Columbia Analytical Services, Inc. | 2655 Park Center Drive, Suite A | Simi Valley, CA 93065 | 805.526.7161 | 805.526.7270 fax | www.caslab.com

LABORATORY REPORT

August 31, 2009

Brian Baker
Environmental Health & Engineering, Inc.
117 Fourth Avenue
Needham, MA 02494

RE: 16512

Dear Brian:

Enclosed are the results of the samples submitted to our laboratory on August 12, 2009. For your reference, these analyses have been assigned our service request number P0902772.

All analyses were performed according to our laboratory's NELAP-approved quality assurance program. The test results meet requirements of the current NELAP standards, where applicable, and except as noted in the laboratory case narrative provided. For a specific list of NELAP-accredited analytes, refer to the certifications section at www.caslab.com. Results are intended to be considered in their entirety and apply only to the samples analyzed and reported herein. Your report contains 526 pages.

Columbia Analytical Services, Inc. is certified by the California Department of Health Services, NELAP Laboratory Certificate No. 02115CA; Arizona Department of Health Services, Certificate No. AZ0694; Florida Department of Health, NELAP Certification E871020; New Jersey Department of Environmental Protection, NELAP Laboratory Certification ID #CA009; New York State Department of Health, NELAP NY Lab ID No: 11221; Oregon Environmental Laboratory Accreditation Program, NELAP ID: CA20007; The American Industrial Hygiene Association, Laboratory #101661; Department of the Navy (NFESC); Pennsylvania Registration No. 68-03307; TX Commission of Environmental Quality, NELAP ID T104704413-08-TX. Each of the certifications listed above have an explicit Scope of Accreditation that applies to specific matrices/methods/analytes; therefore, please contact me for information corresponding to a particular certification.

If you have any questions, please call me at (805) 526-7161.

Respectfully submitted,

Columbia Analytical Services, Inc.

Kate Aguilera

Kate Aguilera
Project Manager

Client: Environmental Health & Engineering, Inc. CAS Project No: P0902772
Project: 16512

CASE NARRATIVE

The samples were received intact under chain of custody on August 12, 2009 and were stored in accordance with the analytical method requirements. Please refer to the sample acceptance check form for additional information. The results reported herein are applicable only to the condition of the samples at the time of sample receipt. The sample labeled "100308" was received wet.

Aldehyde Analysis

The samples were analyzed for aldehydes according to EPA Method TO-11A using high performance liquid chromatography (HPLC).

The results of analyses are given in the attached laboratory report. All results are intended to be considered in their entirety, and Columbia Analytical Services, Inc. (CAS) is not responsible for utilization of less than the complete report.

Client: Environmental Health & Engineering, Incorporate
Project: 16512

Service Request: P0902772

SAMPLE CROSS-REFERENCE

SAMPLE #	CLIENT SAMPLE ID	DATE	TIME
P0902772-001	100303	8/10/09	00:00
P0902772-002	100302	8/10/09	00:00
P0902772-003	100305	8/10/09	00:00
P0902772-004	100306	8/10/09	00:00
P0902772-005	100308	8/10/09	00:00
P0902772-006	100316	8/10/09	00:00
P0902772-007	100443	8/10/09	00:00
P0902772-008	100444	8/10/09	00:00
P0902772-009	100454	8/10/09	00:00
P0902772-010	100451	8/10/09	00:00
P0902772-011	100457	8/10/09	00:00
P0902772-012	100436	8/10/09	00:00

Environmental Health & Engineering, Inc.

CHAIN OF CUSTODY FORM

DATE: 8/10/09

FROM: Environmental Health and Engineering, Inc.
117 Fourth Avenue
Needham, MA 02494-2725

P0902772

TO: Columbia Analytical Svcs

Please send invoices to ATTN: Accounts Payable
Please send reports to ATTN: Data Coordinator

In all correspondence regarding this matter, please refer to EH&E Project # 16512

The cost of this analysis will be covered by EH&E Purchase Order # 16512

For EH & E Data Coordinator - URGENT DATA ☐

	SAMPLE ID	SAMPLE TYPE	ANALYTICAL METHOD/NUMBER	OTHER:Time/Date/Vol.
1	100303	SORBENT TUBES	EPA TO-11	200 MIN
2	100302	↓	↓	↓
3	100305			
4	100306			
5	100308			
6	100316			
7	100443			
8	100444			
9	100454			
10	100451			
11	100457			
12	100436	↓	↓	↓

Special instructions:

☒ Standard turn around time ☐ Rush by _____ (date/time) ☐ Other _____

☐ Fax results 781-247-4305

☐ **RETURN SAMPLES** ☒ Electronic transfer - datacoordinator@eheinc.com

☒ Additional report recipient MFRAGALA@EHEINC.COM

Each signatory please return one copy of this form to the above address

Relinquished by: Margaret Sewell of Environmental Health & Engineering, Inc. Date: 8/11/09

Received by: _____ of (company name) CAS Date: 8/11/09 0945

Relinquished by: _____ of (company name) _____ Date: _____ 20°C

Received by: _____ of (company name) _____ Date: _____

Relinquished by: _____ of (company name) _____ Date: _____

Received by: _____ of (company name) _____ Date: _____

Lab Data
Received by: _____ of Environmental Health & Engineering, Inc. Date: _____

Page ____ of ____

WHITE-EH&E FILE COPY YELLOW-LAB COPY PINK-PROJECT MANAGER COPY GOLD-DATA COORDINATOR COPY

Columbia Analytical Services, Inc.

Sample Acceptance Check Form

Client: Environmental Health & Engineering, Incorporated Work order: P0902772

Project: 16512

Sample(s) received on: 08/12/09 Date opened: 08/12/09 by: MZAMORA

Note: This form is used for all samples received by CAS. The use of this form for custody seals is strictly meant to indicate presence/absence and not as an indication of compliance or nonconformity. Thermal preservation and pH will only be evaluated either at the request of the client and/or as required by the method/SOP.

		Yes	No	N/A
1	Were **sample containers** properly marked with client sample ID?	☒	☐	☐
2	Container(s) **supplied by CAS**?	☒	☐	☐
3	Did **sample containers** arrive in good condition?	☒	☐	☐
4	Was a **chain-of-custody** provided?	☒	☐	☐
5	Was the **chain-of-custody** properly completed?	☐	☒	☐
6	Did **sample container labels** and/or tags agree with custody papers?	☒	☐	☐
7	Was **sample volume** received adequate for analysis?	☒	☐	☐
8	Are samples within specified holding times?	☒	☐	☐
9	Was proper **temperature** (thermal preservation) of cooler at receipt adhered to?	☐	☒	☐
	Cooler Temperature 20 °C Blank Temperature _____ °C			
10	Was a **trip blank** received?	☐	☒	☐
	Trip blank supplied by CAS: _____			
11	Were **custody seals** on outside of cooler/Box?	☐	☒	☐
	Location of seal(s)? _____ Sealing Lid?	☐	☐	☒
	Were signature and date included?	☐	☐	☒
	Were seals intact?	☐	☐	☒
	Were custody seals on outside of sample container?	☐	☒	☐
	Location of seal(s)? _____ Sealing Lid?	☐	☐	☒
	Were signature and date included?	☐	☐	☒
	Were seals intact?	☐	☐	☒
12	Do containers have appropriate **preservation**, according to method/SOP or Client specified information?	☐	☐	☒
	Is there a client indication that the submitted samples are **pH** preserved?	☐	☐	☒
	Were **VOA vials** checked for presence/absence of air bubbles?	☐	☐	☒
	Does the client/method/SOP require that the analyst check the sample pH and if necessary alter it?	☐	☐	☒
13	**Tubes:** Are the tubes capped and intact?	☒	☐	☐
	Do they contain moisture?	☐	☒	☐
14	**Badges:** Are the badges properly capped and intact?	☐	☐	☒
	Are dual bed badges separated and individually capped and intact?	☐	☐	☒

Lab Sample ID	Container Description	Required pH *	Received pH	Adjusted pH	VOA Headspace (Presence/Absence)	Receipt / Preservation Comments
P0902772-001.01	Silica Gel DNPH Tube					
P0902772-002.01	Silica Gel DNPH Tube					
P0902772-003.01	Silica Gel DNPH Tube					
P0902772-004.01	Silica Gel DNPH Tube					
P0902772-005.01	Silica Gel DNPH Tube					
P0902772-006.01	Silica Gel DNPH Tube					

Explain any discrepancies: (include lab sample ID numbers):

Chain of Custody is missing time collected

Sampling time is listed on COC (200 min),flow rate and total volume are not listed.

*Required pH: Phenols/COD/NH3/TOC/TOX/NO3+NO2/TKN/T.PHOS, H2SO4 (pH<2); Metals, HNO3 (pH<2); CN (NaOH or NaOH/Asc Acid) (pH>12); Diss. Sulfide, NaOH (pH>12); T. Sulfide, NaOH/ZnAc (pH>12) RSK - MEEPP, HCL (pH<2); RSK - CO2, (pH 5-8); Sulfur (pH>4)

Columbia Analytical Services, Inc.
Sample Acceptance Check Form

Client: Environmental Health & Engineering, Incorporated Work order: P0902772

Project: 16512

Sample(s) received on: 08/12/09 Date opened: 08/12/09 by: MZAMORA

Lab Sample ID	Container Description	Required pH *	Received pH	Adjusted pH	VOA Headspace (Presence/Absence)	Receipt / Preservation Comments
P0902772-007.01	Silica Gel DNPH Tube					
P0902772-008.01	Silica Gel DNPH Tube					
P0902772-009.01	Silica Gel DNPH Tube					
P0902772-010.01	Silica Gel DNPH Tube					
P0902772-011.01	Silica Gel DNPH Tube					
P0902772-012.01	Silica Gel DNPH Tube					

Explain any discrepancies: (include lab sample ID numbers):

*Required pH: Phenols/COD/NH3/TOC/TOX/NO3+NO2/TKN/T.PHOS, H2SO4 (pH<2); Metals, HNO3 (pH<2); CN (NaOH or NaOH/Asc Acid) (pH>12); Diss. Sulfide, NaOH (pH>12); T. Sulfide, NaOH/ZnAc (pH>12) RSK - MEEPP, HCL (pH<2); RSK - CO2, (pH 5-8); Sulfur (pH>4)

RESULTS OF ANALYSIS

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **100303**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P0902772-001

Test Code: EPA Method TO-11A
Instrument ID: Waters LC Module I Plus/UV_Vis 360/LC1
Analyst: Hani Cherazaie
Sampling Media: Silica Gel DNPH Tube
Test Notes: **BC**

Date Collected: 8/10/09
Date Received: 8/12/09
Date Analyzed: 8/17 - 8/18/09
Desorption Volume: 1.0 ml
Volume Sampled: NA Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	Formaldehyde	< 100	NA	NA	NA	NA	
75-07-0	Acetaldehyde	< 100	NA	NA	NA	NA	
123-38-6	Propionaldehyde	< 100	NA	NA	NA	NA	
4170-30-3	Crotonaldehyde, Total	< 100	NA	NA	NA	NA	
123-72-8	Butyraldehyde	< 100	NA	NA	NA	NA	
100-52-7	Benzaldehyde	< 100	NA	NA	NA	NA	
590-86-3	Isovaleraldehyde	< 100	NA	NA	NA	NA	
110-62-3	Valeraldehyde	< 100	NA	NA	NA	NA	
529-20-4	o-Tolualdehyde	< 100	NA	NA	NA	NA	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	NA	NA	NA	NA	
66-25-1	n-Hexaldehyde	< 100	NA	NA	NA	NA	
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	NA	NA	NA	NA	

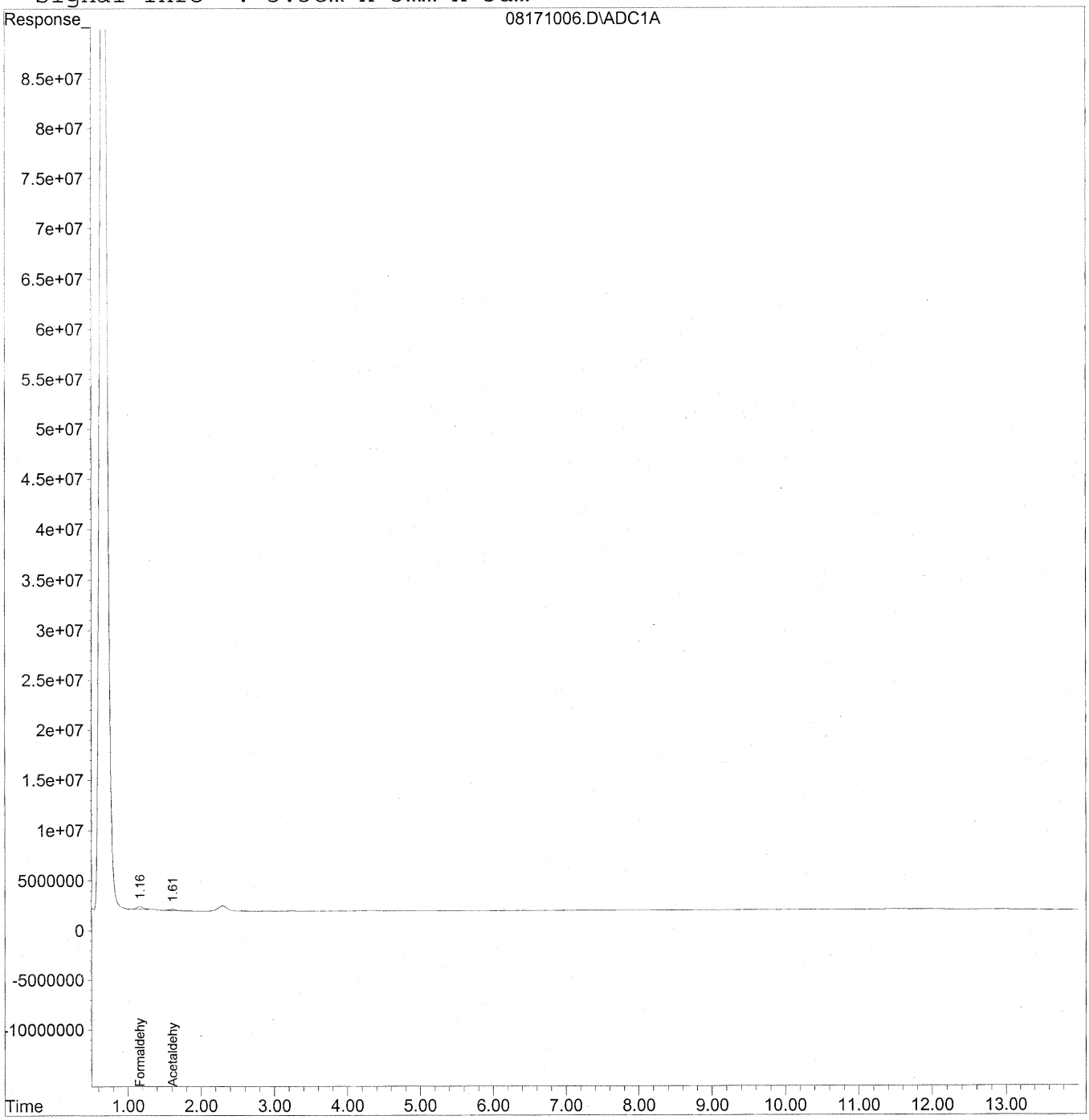
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.

Verified By: _____ Date: 8/25/09

Data File : J:\LC01\DATA\TO11\2009_08\17\08171006.D Vial: 19
Acq On : 18 Aug 2009 5:09 pm Operator: HC
Sample : P0902772-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



9

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171006.D Vial: 19
Acq On : 18 Aug 2009 5:09 pm Operator: HC
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Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

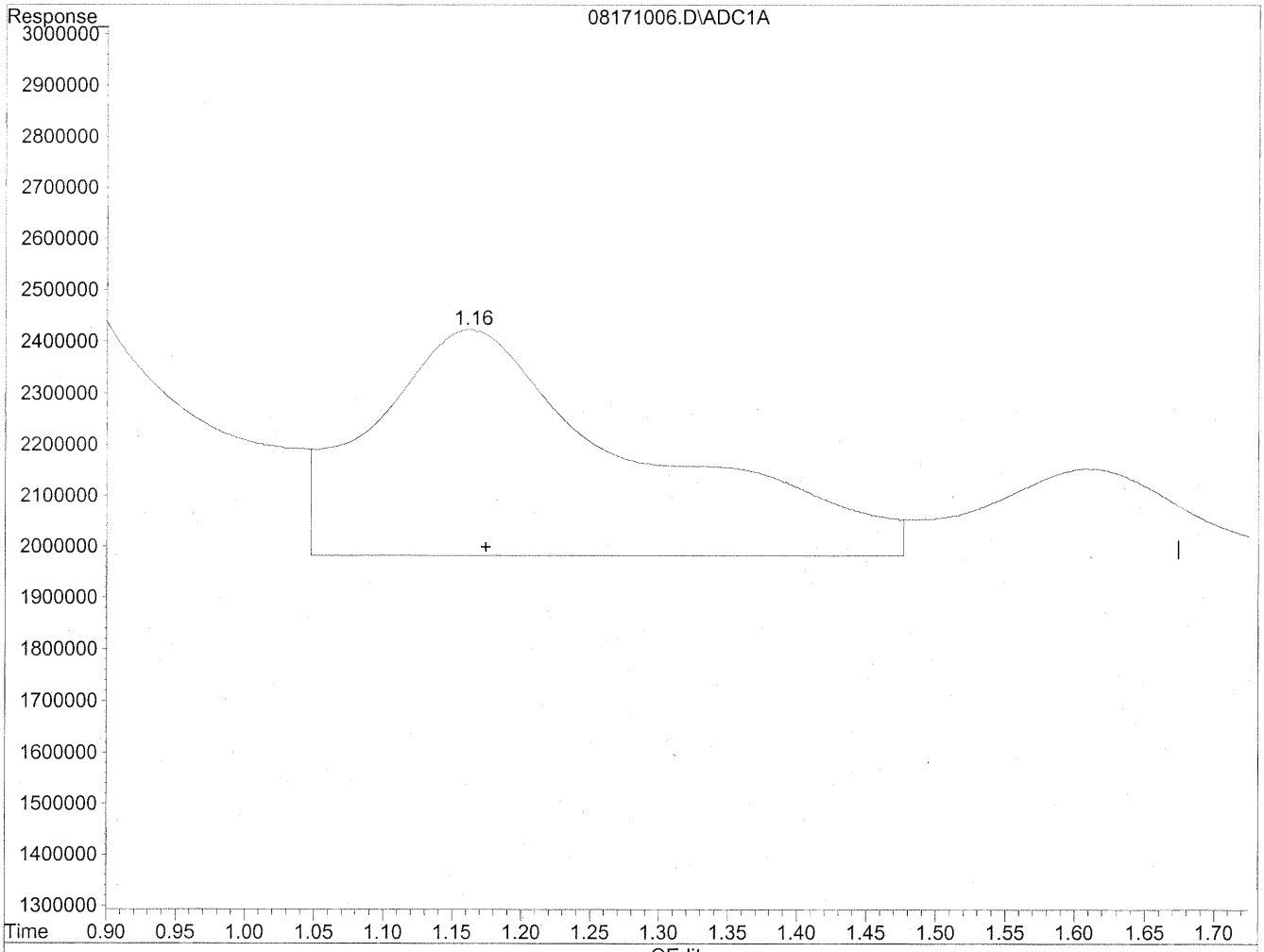
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	15882916	86.517	ng/mlm
2)	Acetaldehyde	1.61	9176088	65.439	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

10

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171006.D Vial: 19
Acq On : 18 Aug 2009 5:09 pm Operator: HC
Sample : P0902772-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.16min 322.941ng/ml

response 59285887

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171006.D Vial: 19
Acq On : 18 Aug 2009 5:09 pm Operator: HC
Sample : P0902772-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

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Title : TO-11A Method for Aldehydes/Ketones by HPLC
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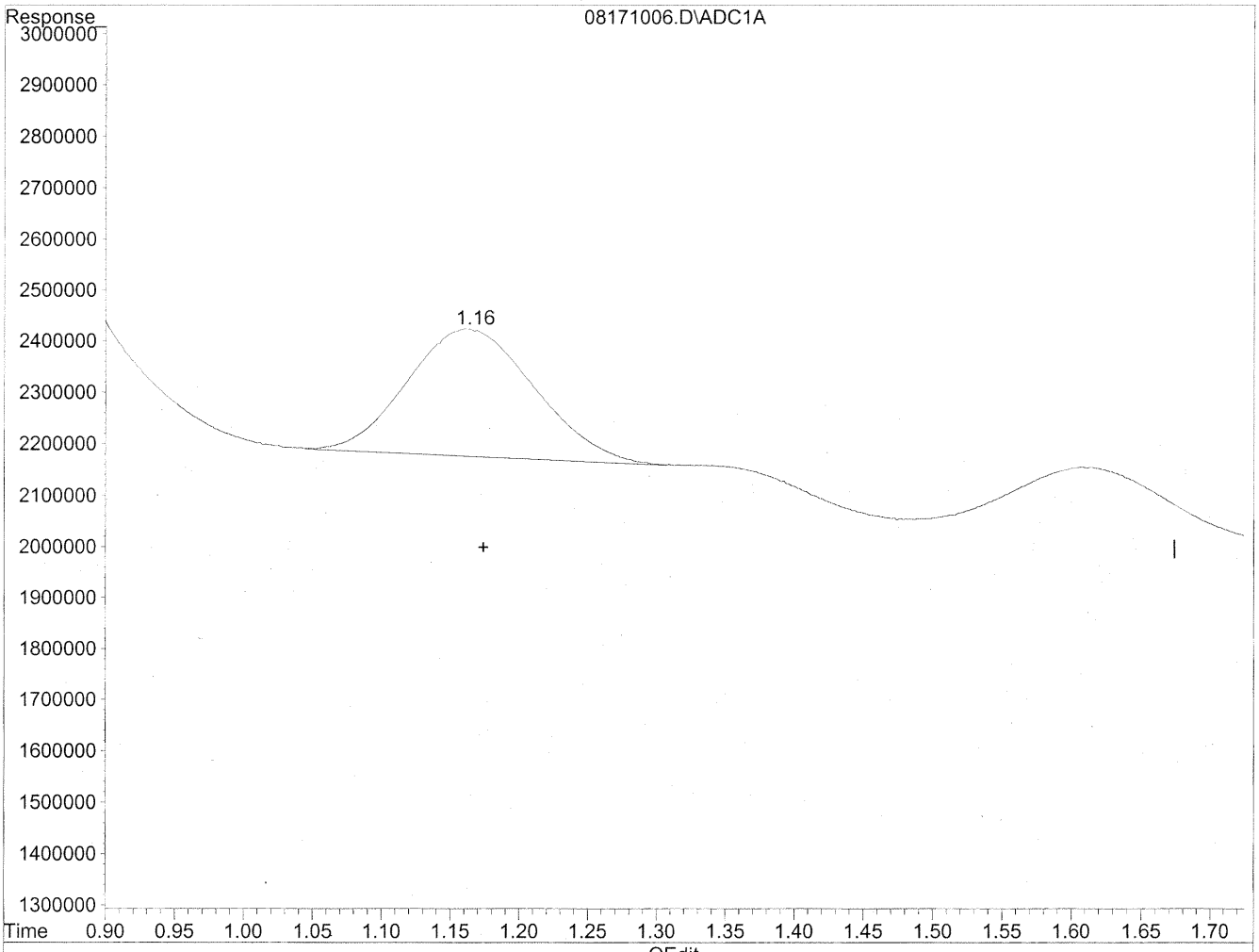


(1) Formaldehyde

1.16min 86.517ng/ml m

response 15882916

HL 8/22/09 LC

KP 8/23/09

12

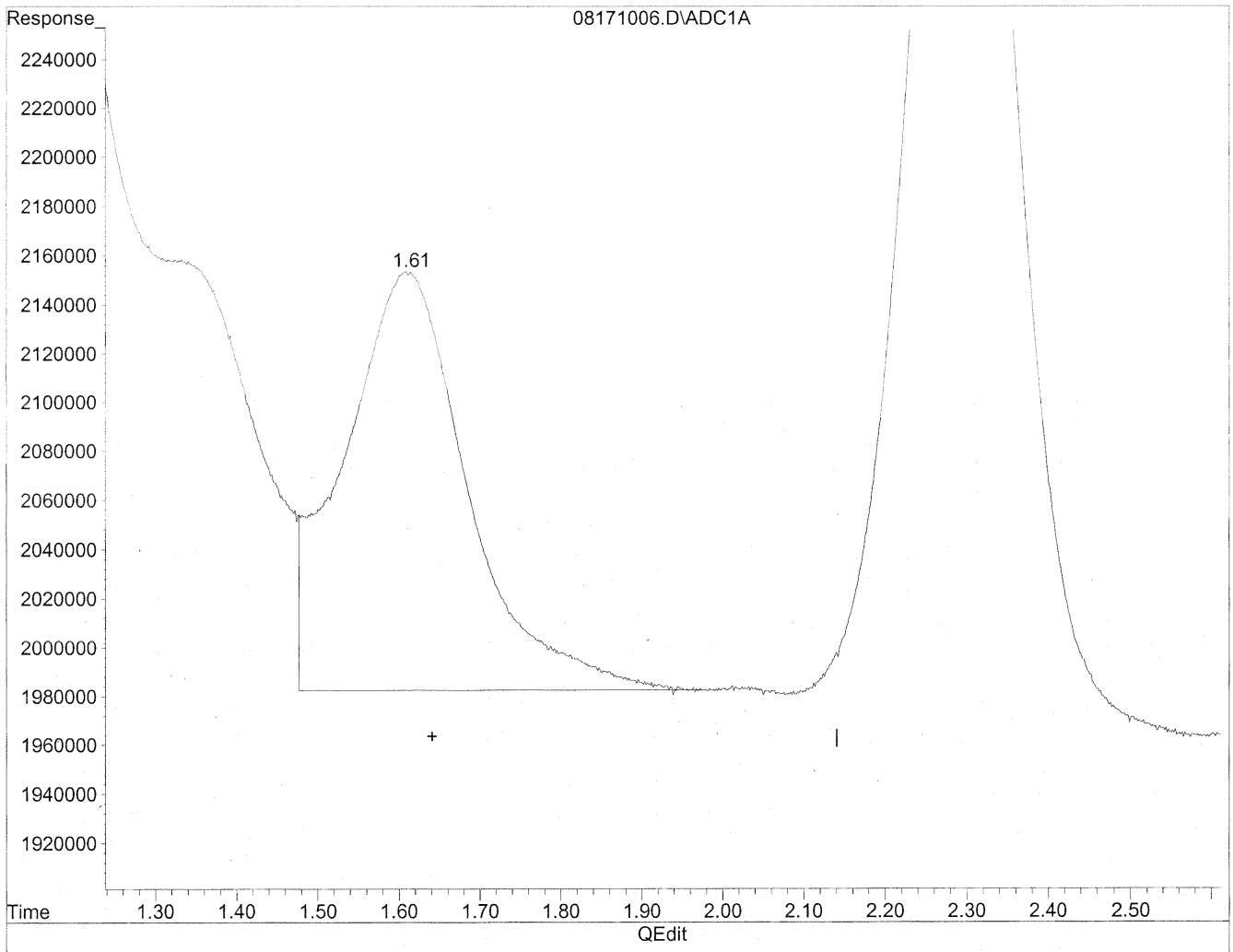
(+) = Expected Retention Time

08171006.D TO110709.M Sat Aug 22 14:08:25 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171006.D Vial: 19
Acq On : 18 Aug 2009 5:09 pm Operator: HC
Sample : P0902772-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.61min 129.578ng/ml

response 18169909

(+) = Expected Retention Time

08171006.D TO110709.M Sat Aug 22 14:08:33 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171006.D Vial: 19
Acq On : 18 Aug 2009 5:09 pm Operator: HC
Sample : P0902772-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
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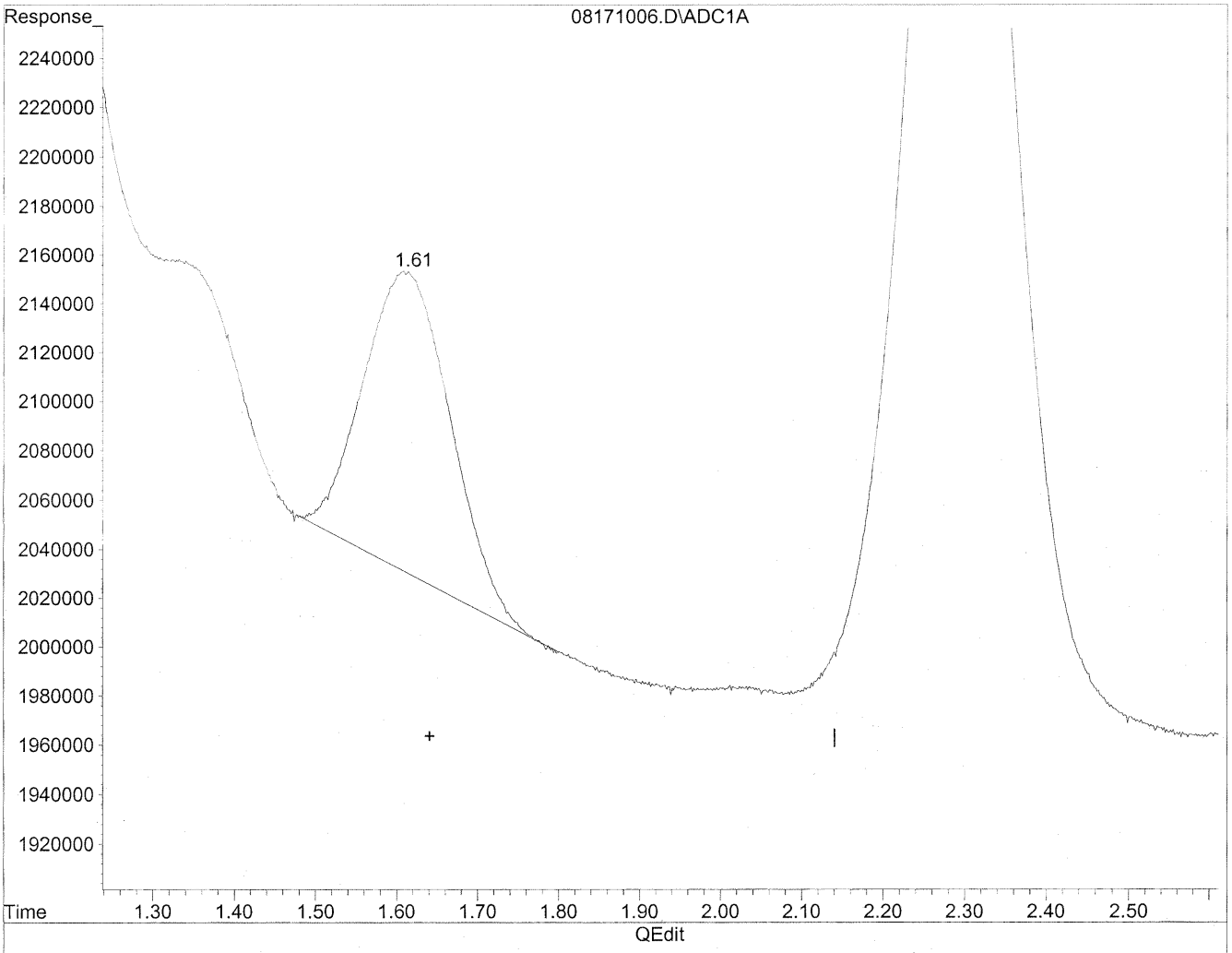


(2) Acetaldehyde

1.61min 65.439ng/ml m

response 9176088

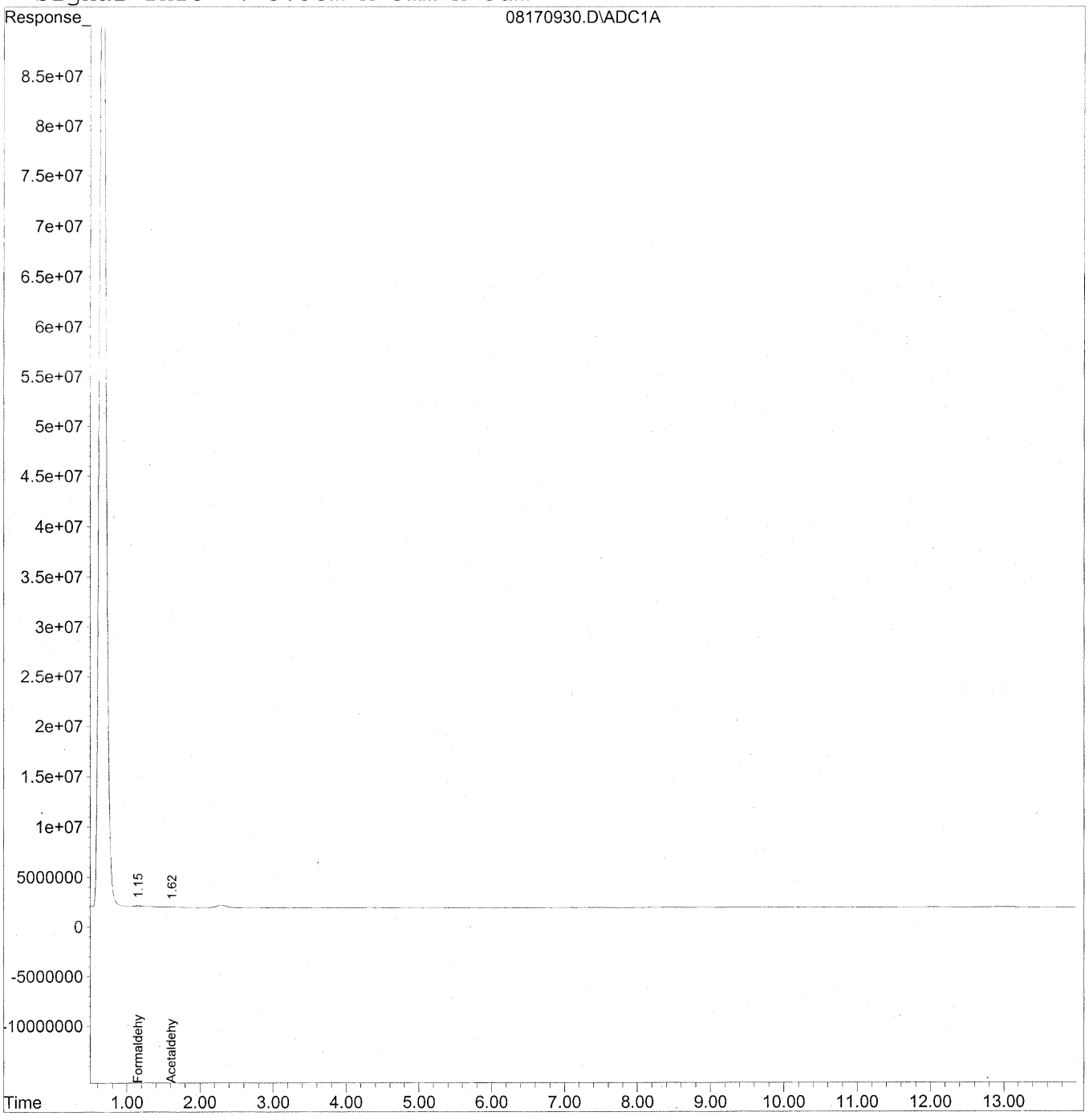
HC 8/22/09
LC
KP 8/23/09

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_08\17\08170930.D Vial: 29
Acq On : 17 Aug 2009 10:06 pm Operator: HC
Sample : P0902772-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:06 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



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Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170930.D Vial: 29
Acq On : 17 Aug 2009 10:06 pm Operator: HC
Sample : P0902772-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:06 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

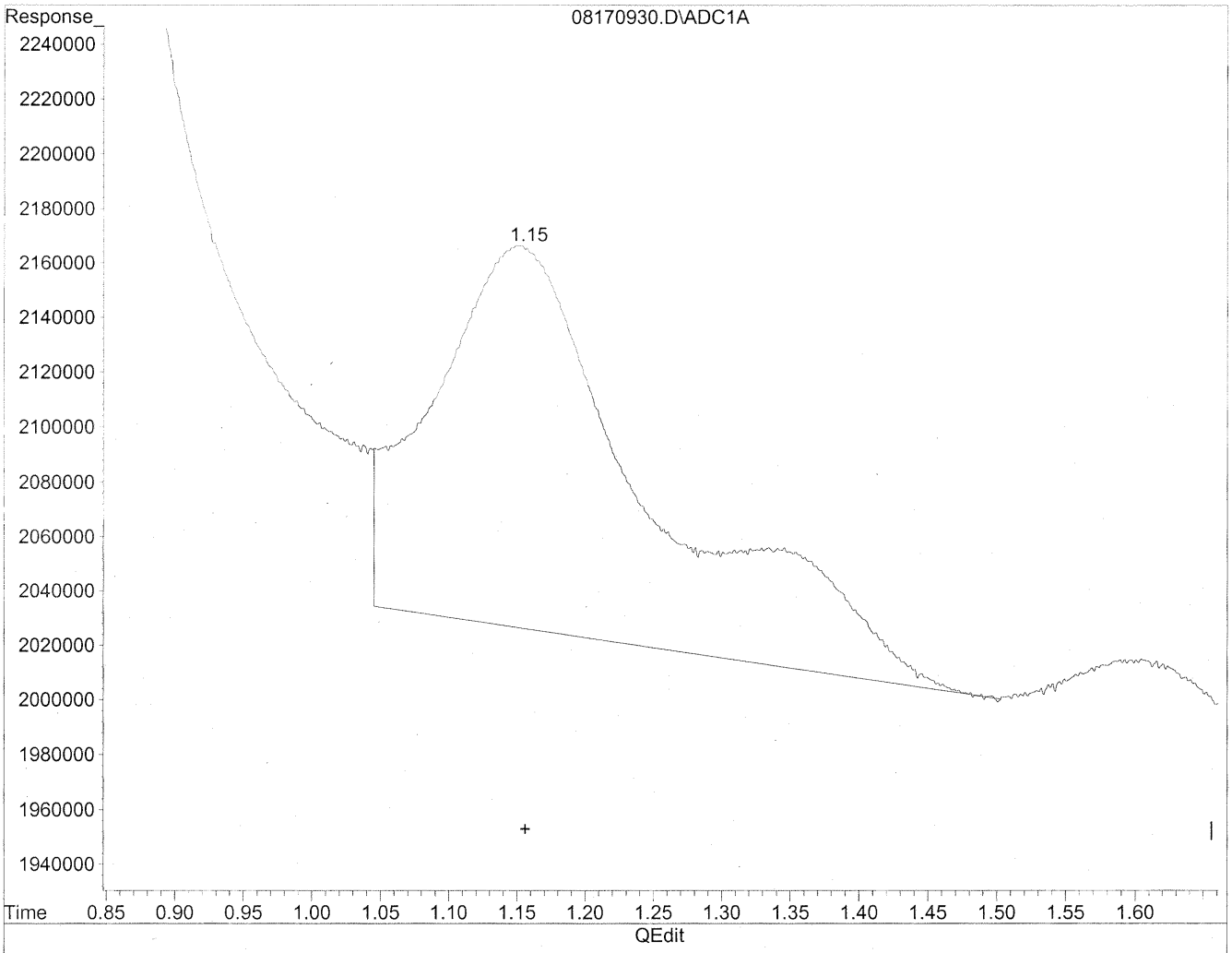
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.15	5619238	30.609	ng/mlm
2)	Acetaldehyde	1.62	1644774	11.730	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170930.D Vial: 29
Acq On : 17 Aug 2009 10:06 pm Operator: HC
Sample : P0902772-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:04 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.15min 84.284ng/ml

response 15472976

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170930.D Vial: 29
Acq On : 17 Aug 2009 10:06 pm Operator: HC
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Misc : Multiplr: 1.00
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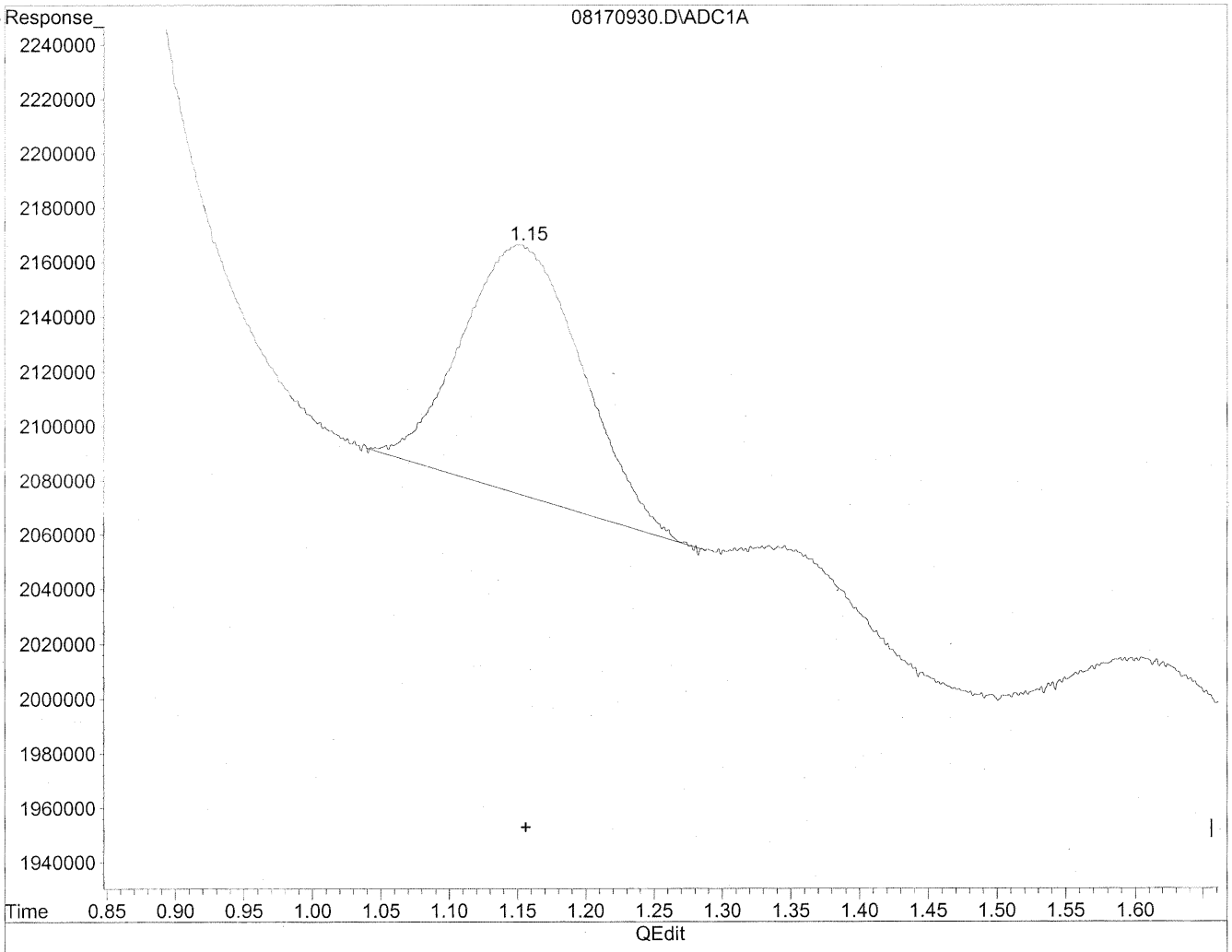

(1) Formaldehyde

1.15min 30.609ng/ml m

response 5619238

HC 8/21/09
IC
KP 8/23/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170930.D Vial: 29
Acq On : 17 Aug 2009 10:06 pm Operator: HC
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Misc : Multiplr: 1.00
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Quant Time: Aug 21 18:04 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
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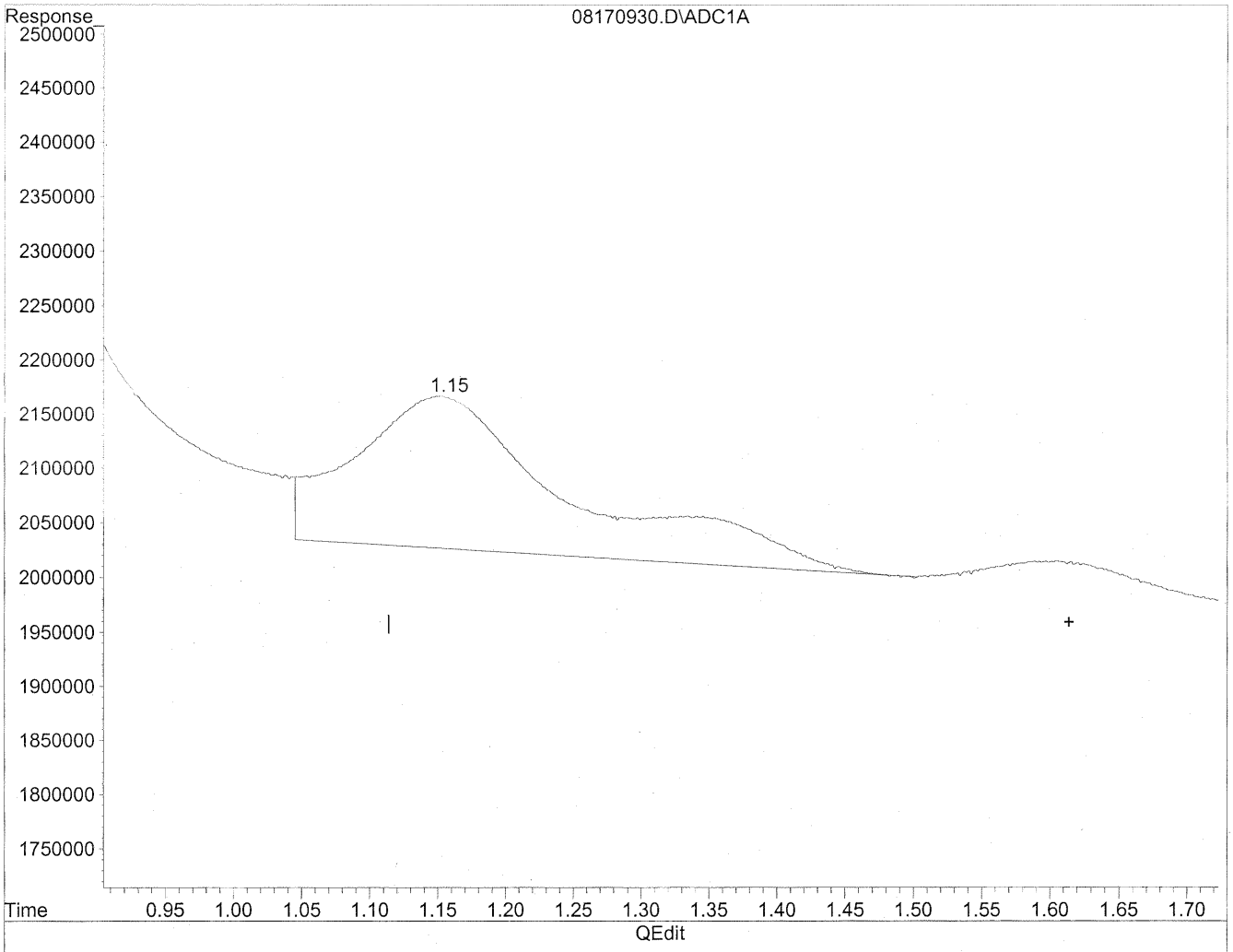


(2) Acetaldehyde

1.15min 110.345ng/ml

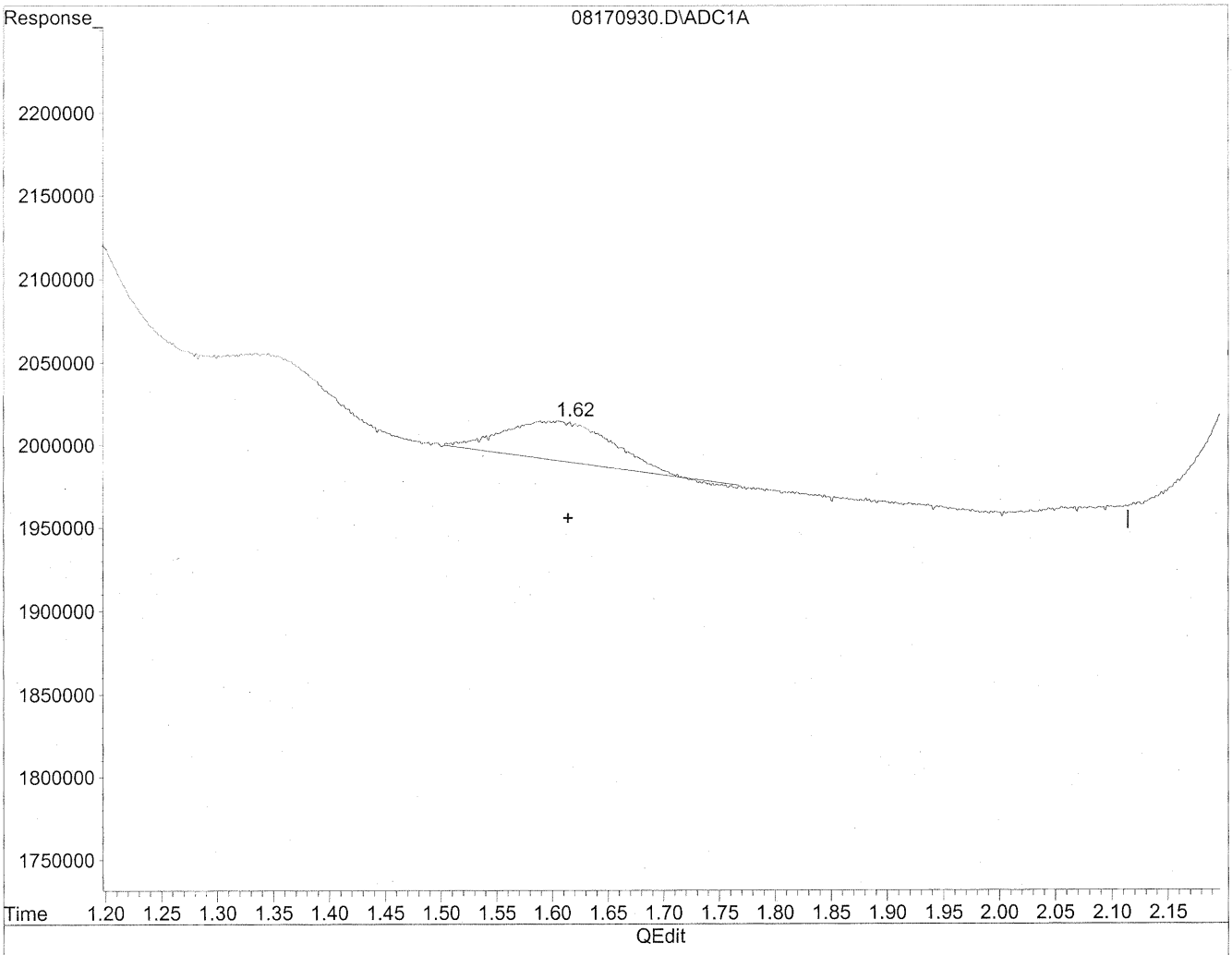
response 15472976

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170930.D Vial: 29
Acq On : 17 Aug 2009 10:06 pm Operator: HC
Sample : P0902772-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:04 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.62min 11.730ng/ml m

response 1644774

HC 8/21/09 MP

KR 8/23/09

20

(+) = Expected Retention Time

08170930.D TO110709.M Fri Aug 21 18:06:17 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **100302**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P0902772-002

Test Code:	EPA Method TO-11A	Date Collected:	8/10/09
Instrument ID:	Waters LC Module I Plus/UV_Vis 360/LC1	Date Received:	8/12/09
Analyst:	Hani Cherazaie	Date Analyzed:	8/17 - 8/18/09
Sampling Media:	Silica Gel DNPH Tube	Desorption Volume:	1.0 ml
Test Notes:	**BC**	Volume Sampled:	105.6 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**3,800**	**36**	0.95	**30**	0.77	
75-07-0	**Acetaldehyde**	**2,600**	**24**	0.95	**14**	0.53	**BT**
123-38-6	**Propionaldehyde**	**200**	**1.9**	0.95	**0.79**	0.40	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.95	ND	0.33	
123-72-8	**Butyraldehyde**	**210**	**2.0**	0.95	**0.67**	0.32	
100-52-7	**Benzaldehyde**	**310**	**3.0**	0.95	**0.68**	0.22	
590-86-3	Isovaleraldehyde	< 100	ND	0.95	ND	0.27	
110-62-3	**Valeraldehyde**	**250**	**2.3**	0.95	**0.66**	0.27	
529-20-4	o-Tolualdehyde	< 100	ND	0.95	ND	0.19	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	1.9	ND	0.39	
66-25-1	**n-Hexaldehyde**	**610**	**5.8**	0.95	**1.4**	0.23	
5779-94-2	**2,5-Dimethylbenzaldehyde**	**490**	**4.6**	0.95	**0.84**	0.17	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
BT = Results indicated possible breakthrough; back section > 10% front section.

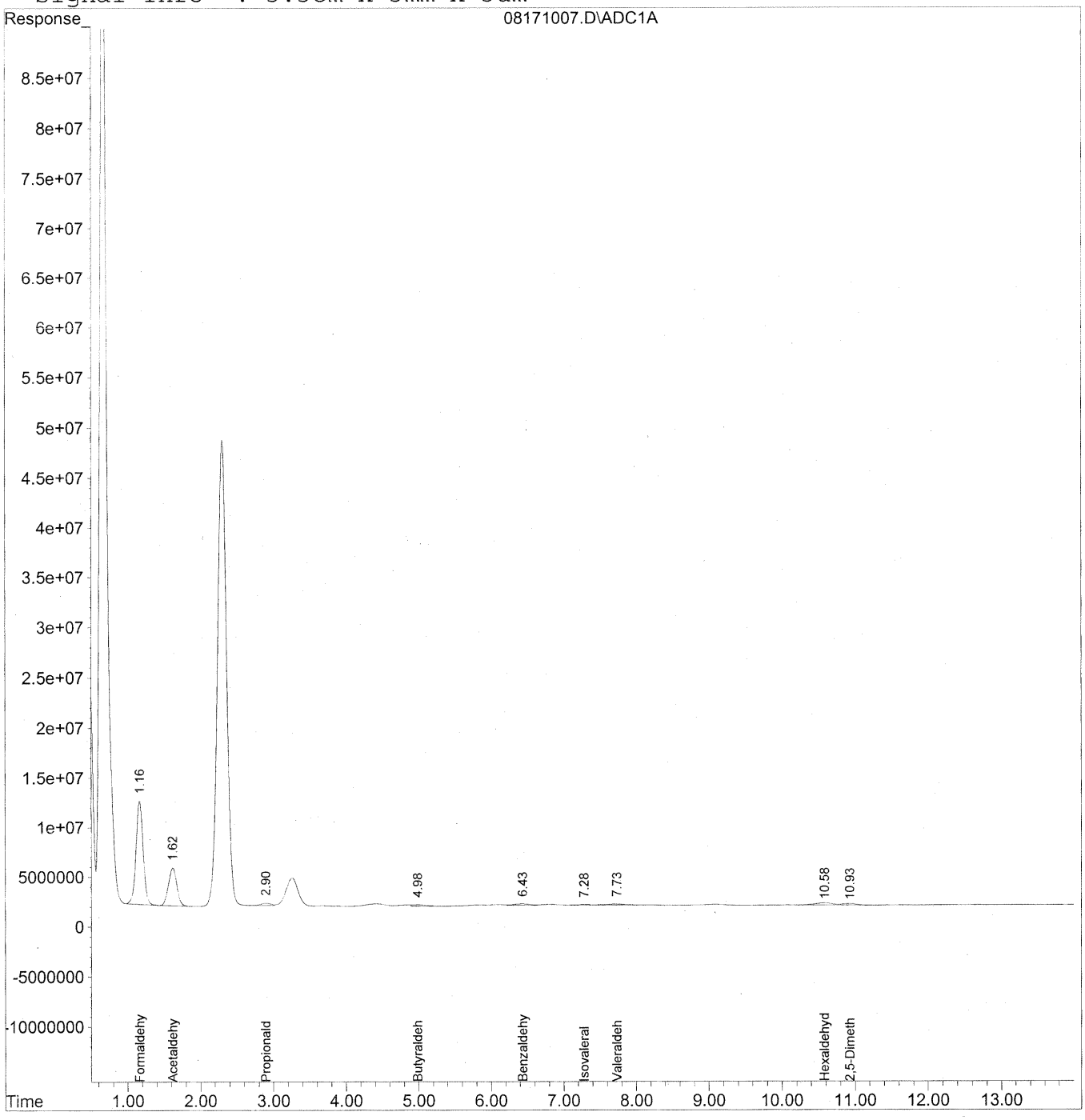
Verified By: _____ Date: 8/25/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



22

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

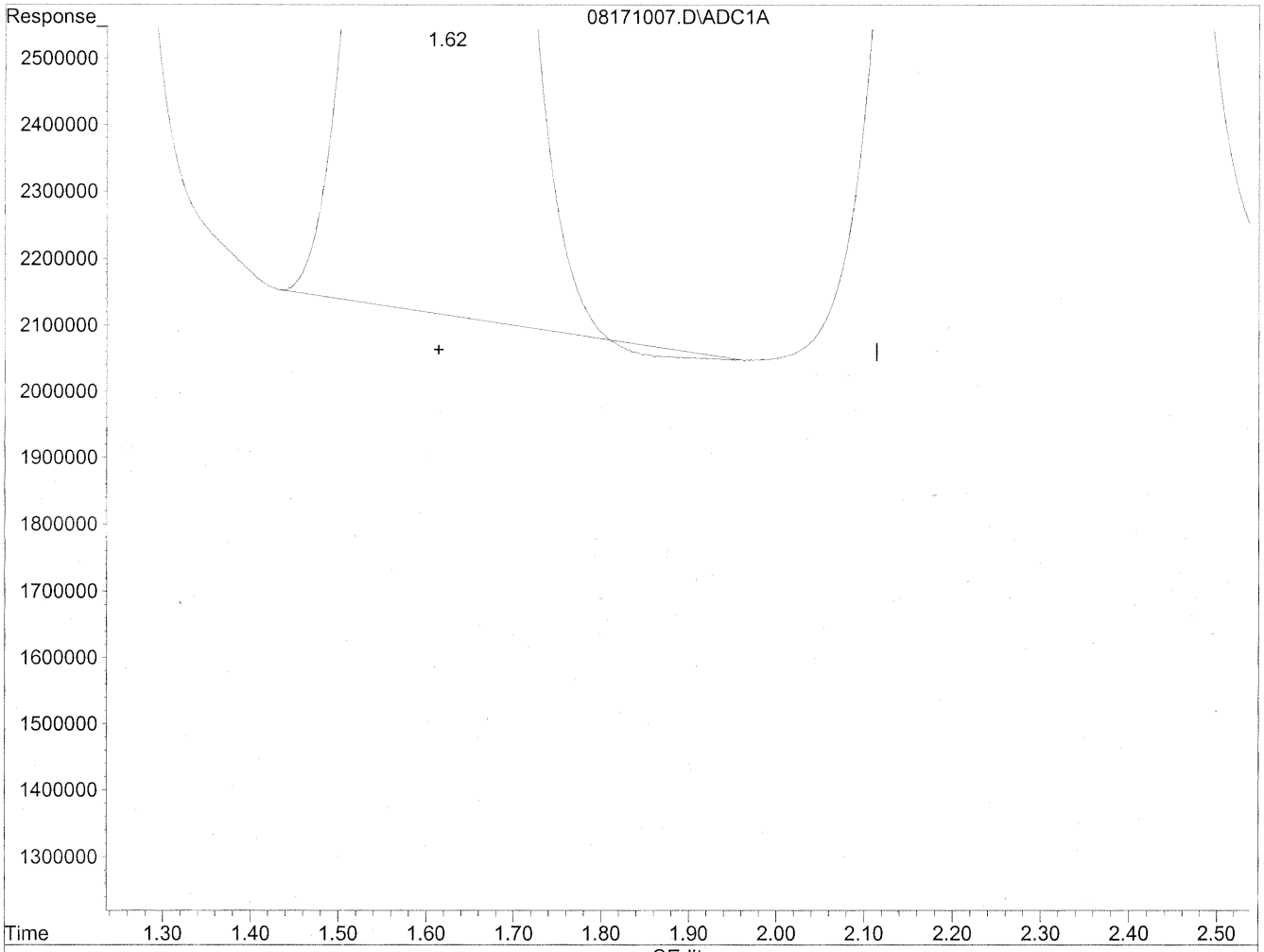
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	705386835	3842.366	ng/ml
2)	Acetaldehyde	1.62	305894827	2181.480	ng/mlm
3)	Propionaldehyde	2.90	21069913	197.478	ng/mlm
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	4.98	18540713	209.888	ng/mlm
6)	Benzaldehyde	6.43	20644571	313.417	ng/mlm
7)	Isovaleraldehyde	7.28	6062965	77.481	ng/mlm
8)	Valeraldehyde	7.73	18076141	245.917	ng/mlm
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/mld
11)	Hexaldehyde	10.58	41277424	612.936	ng/mlm
12)	2,5-Dimethylbenzaldehyde	10.92	23978415	489.222	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.62min 2164.324ng/ml

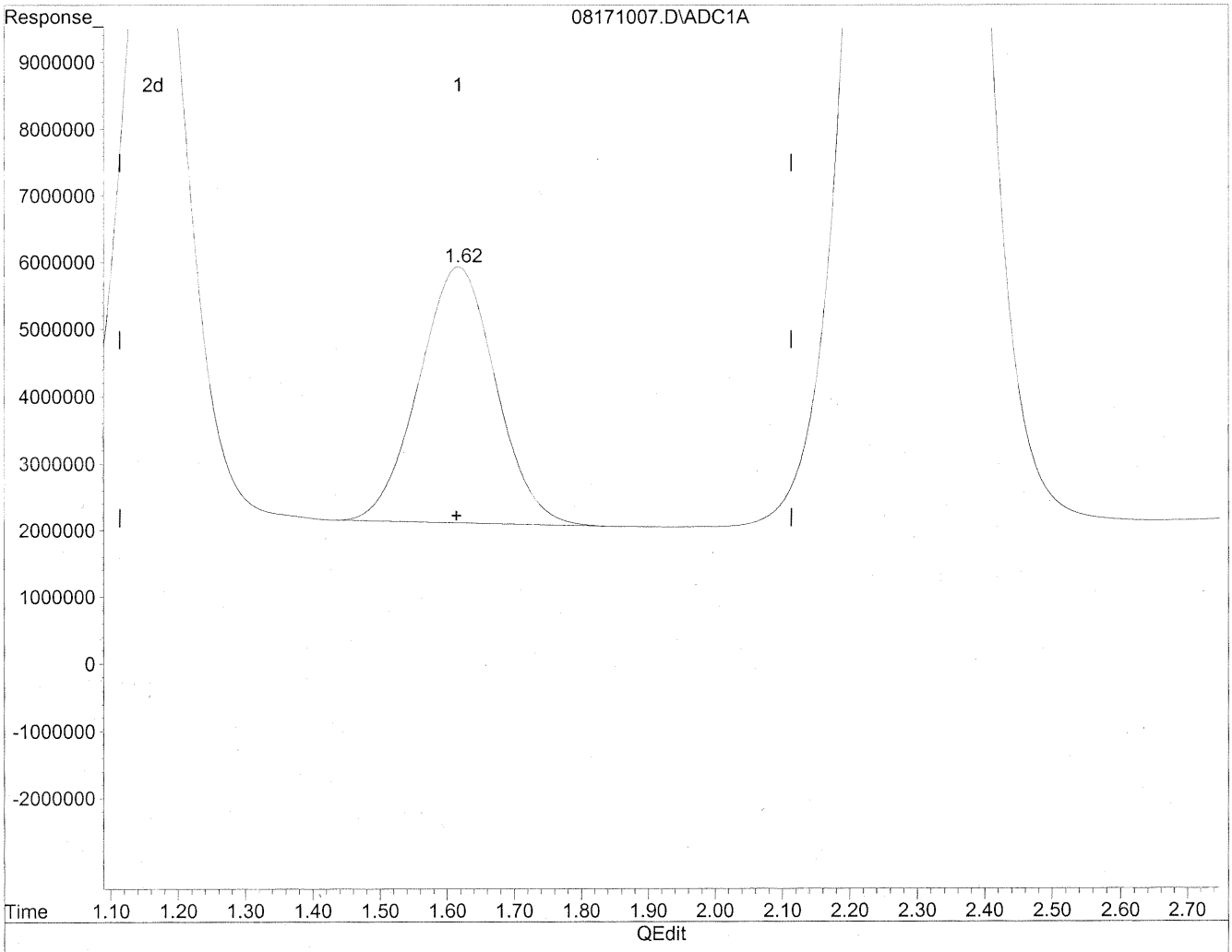
response 303489059

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.62min 2181.480ng/ml m

response 305894827

HC 8/22/09
LC
KR 8/23/09

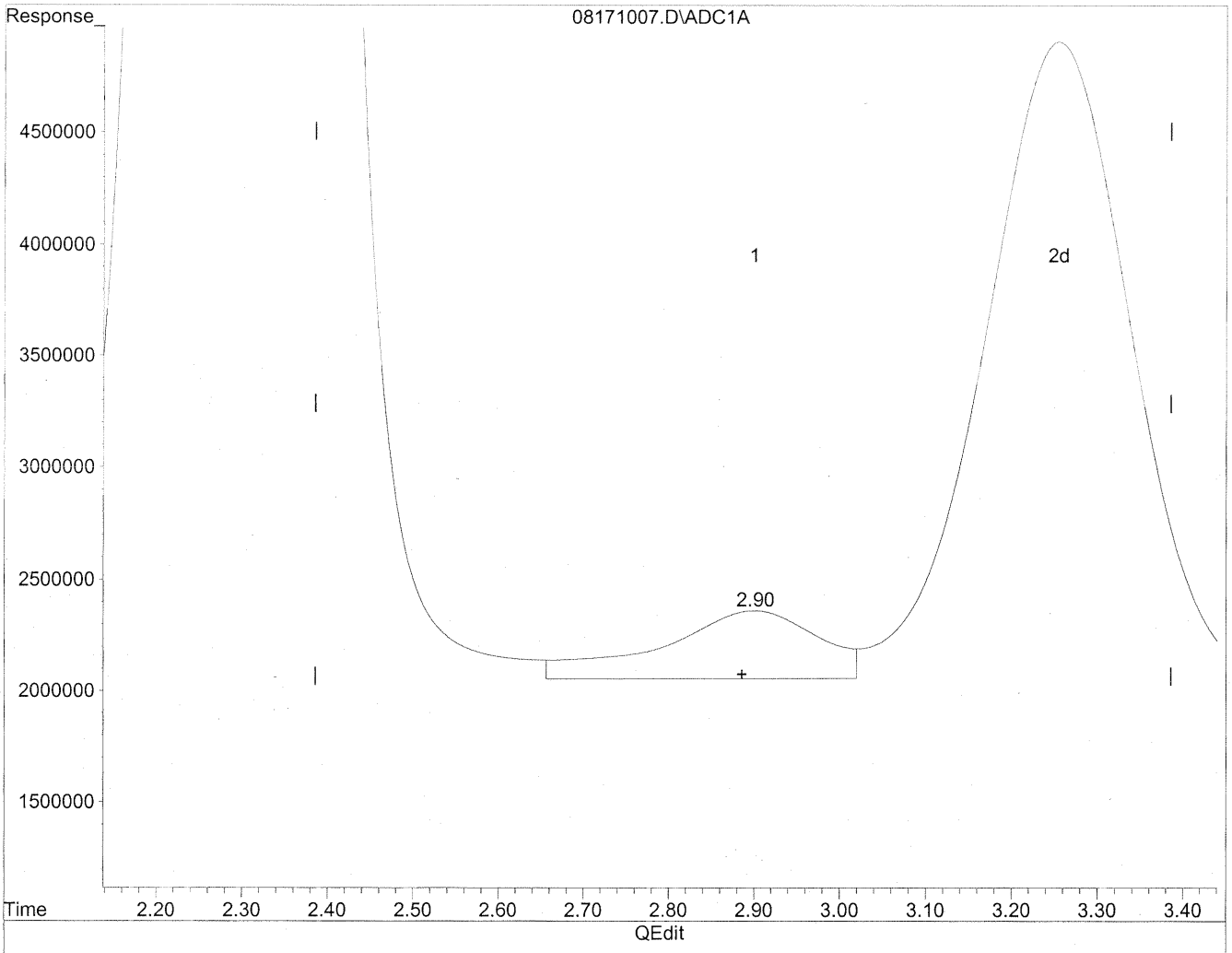
25

(+) = Expected Retention Time
08171007.D TO110709.M Sat Aug 22 14:09:06 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.90min 367.875ng/ml

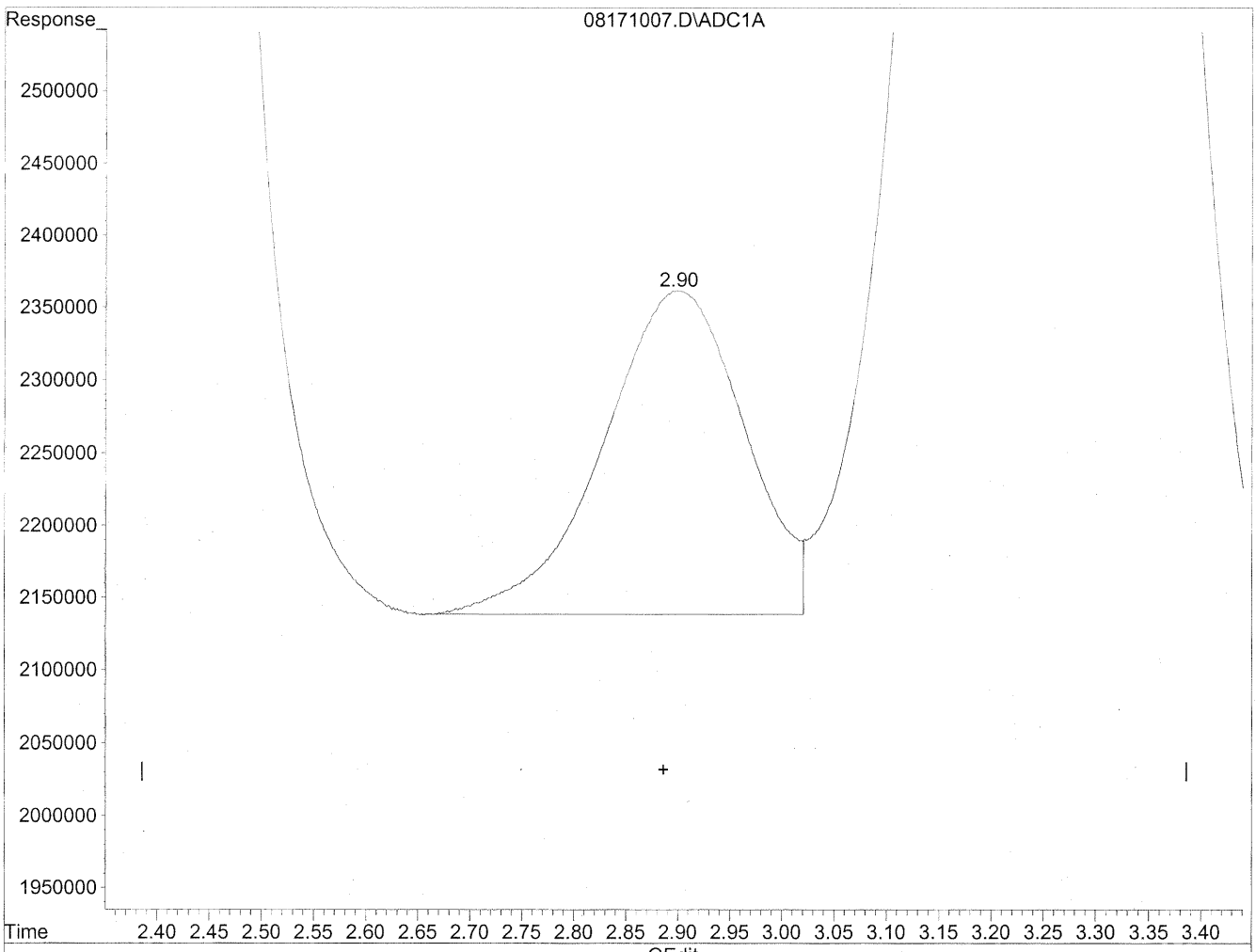
response 39250515

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.90min 197.478ng/ml m

response 21069913

HC 8/22/09
IC
KP 8/23/09

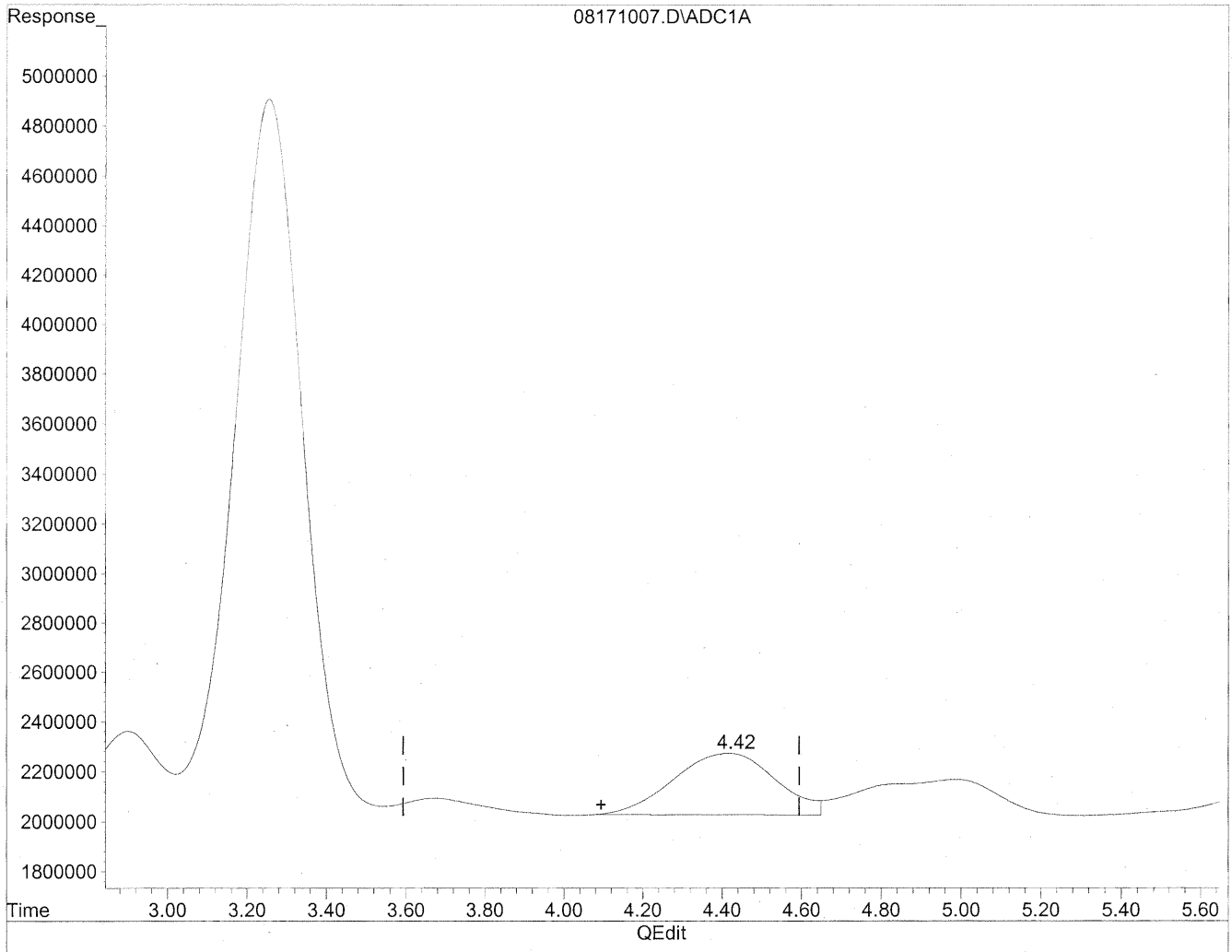
(+) = Expected Retention Time

08171007.D TO110709.M Sat Aug 22 14:09:40 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.42min 451.752ng/ml

response 44007519

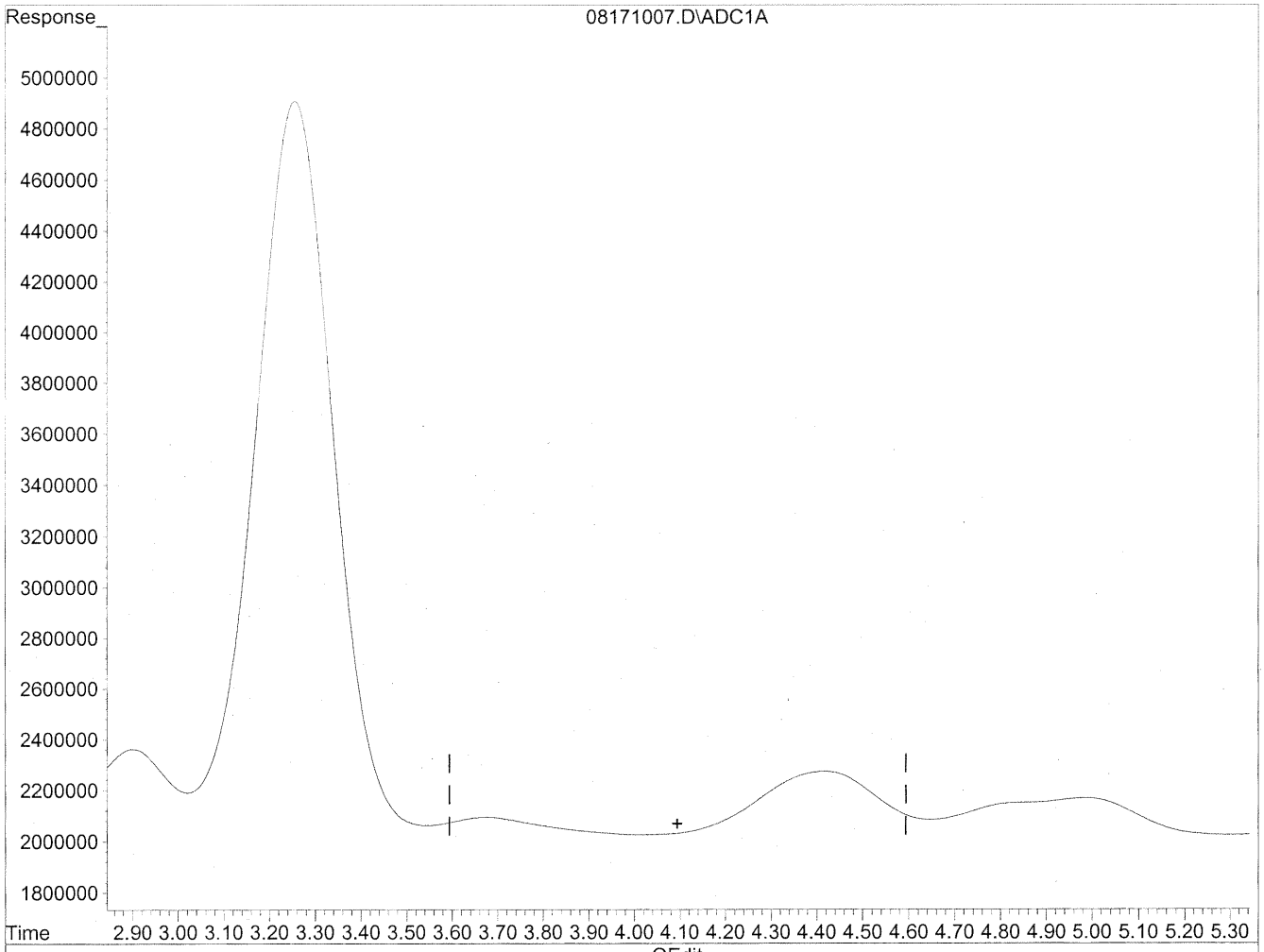
(+) = Expected Retention Time

08171007.D TO110709.M Sat Aug 22 14:09:47 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09 MP

KH 8/23/09

29

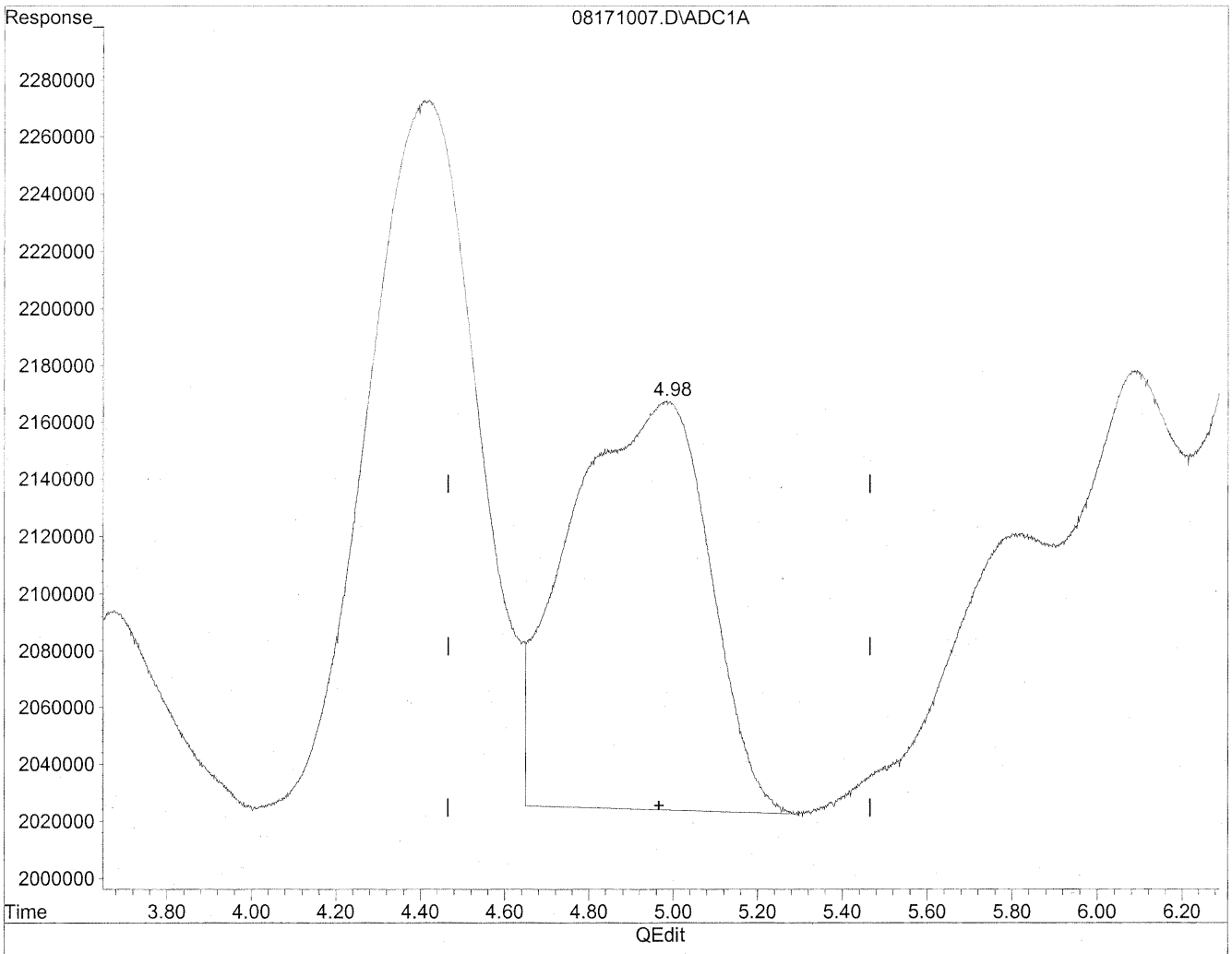
(+) = Expected Retention Time

08171007.D TO110709.M Sat Aug 22 14:09:51 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.98min 372.207ng/ml

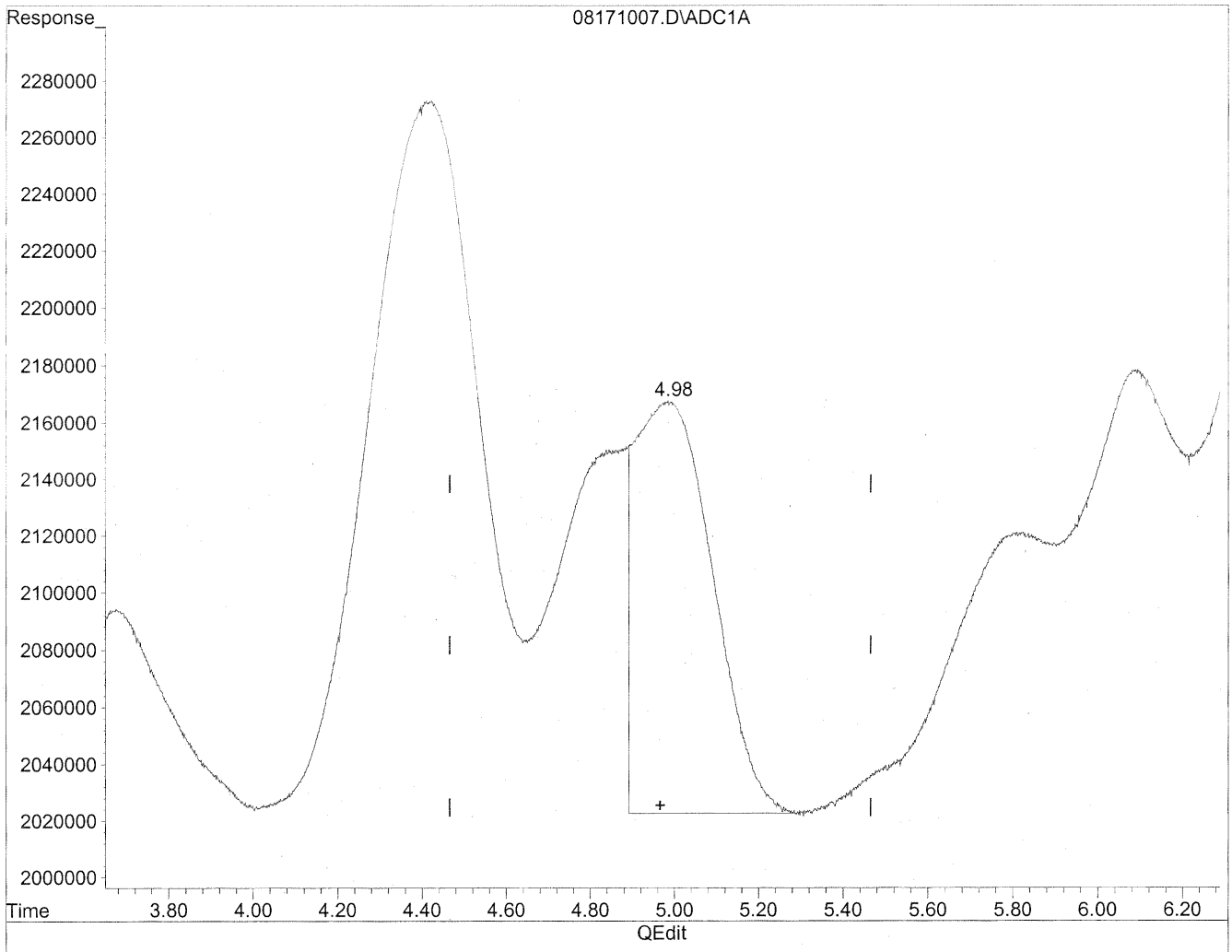
response 32879327

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.98min 209.888ng/ml m

response 18540713

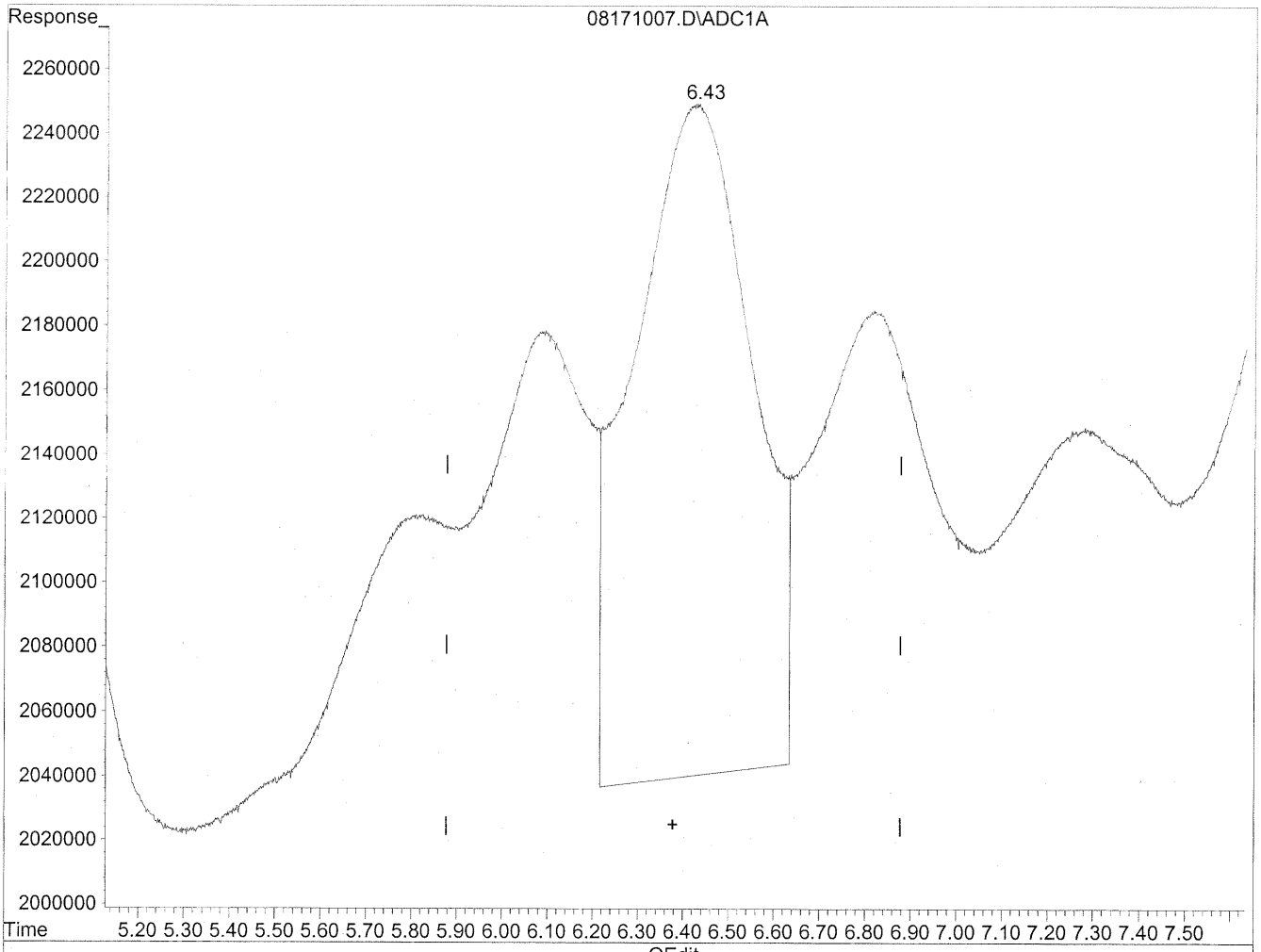
HC 8/22/09
SP
KK 8/23/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(6) Benzaldehyde

6.42min 576.657ng/ml

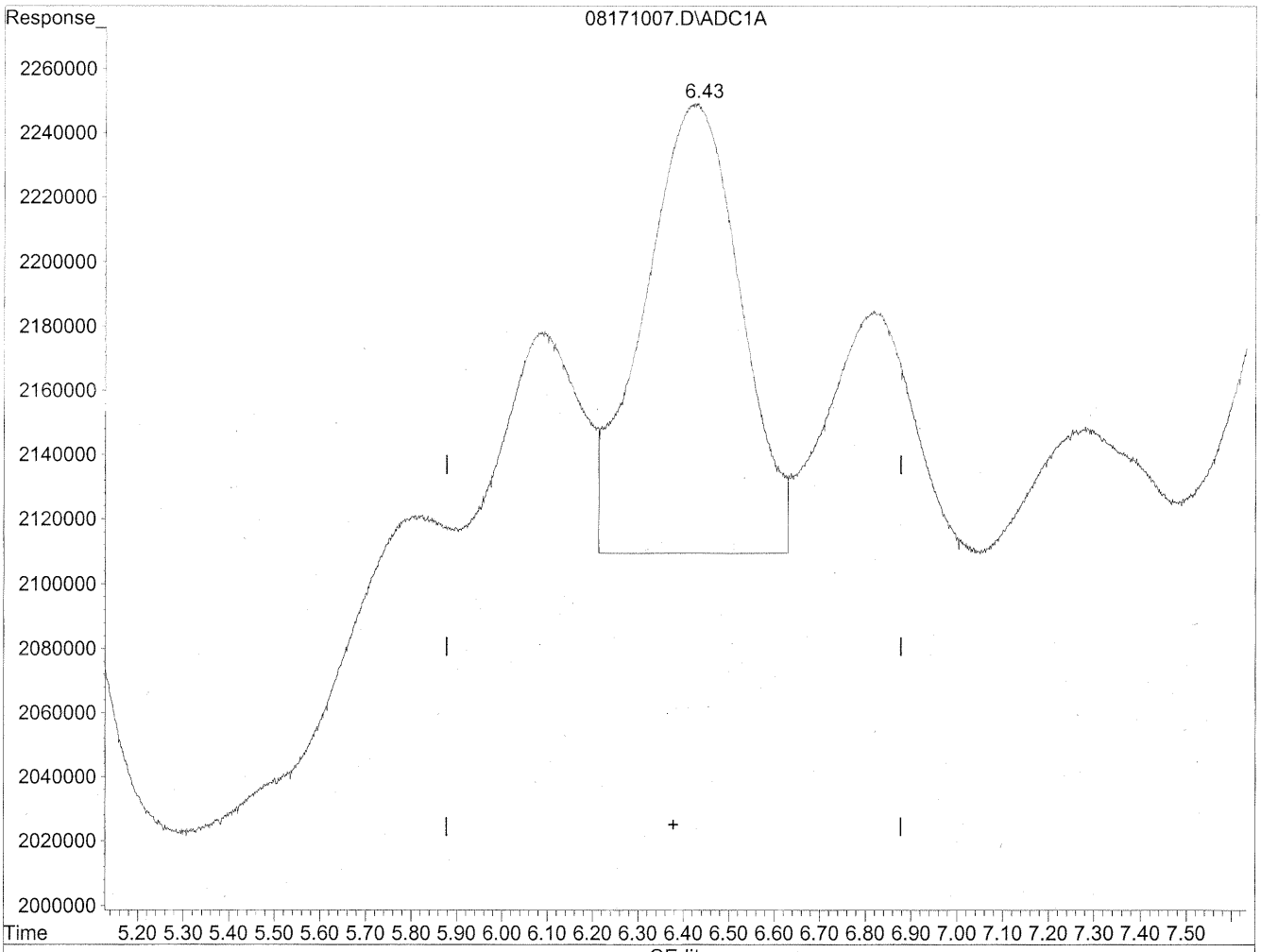
response 37983979

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(6) Benzaldehyde

6.43min 313.417ng/ml m

response 20644571

HC 8/22/09
BC
KP 8/23/09

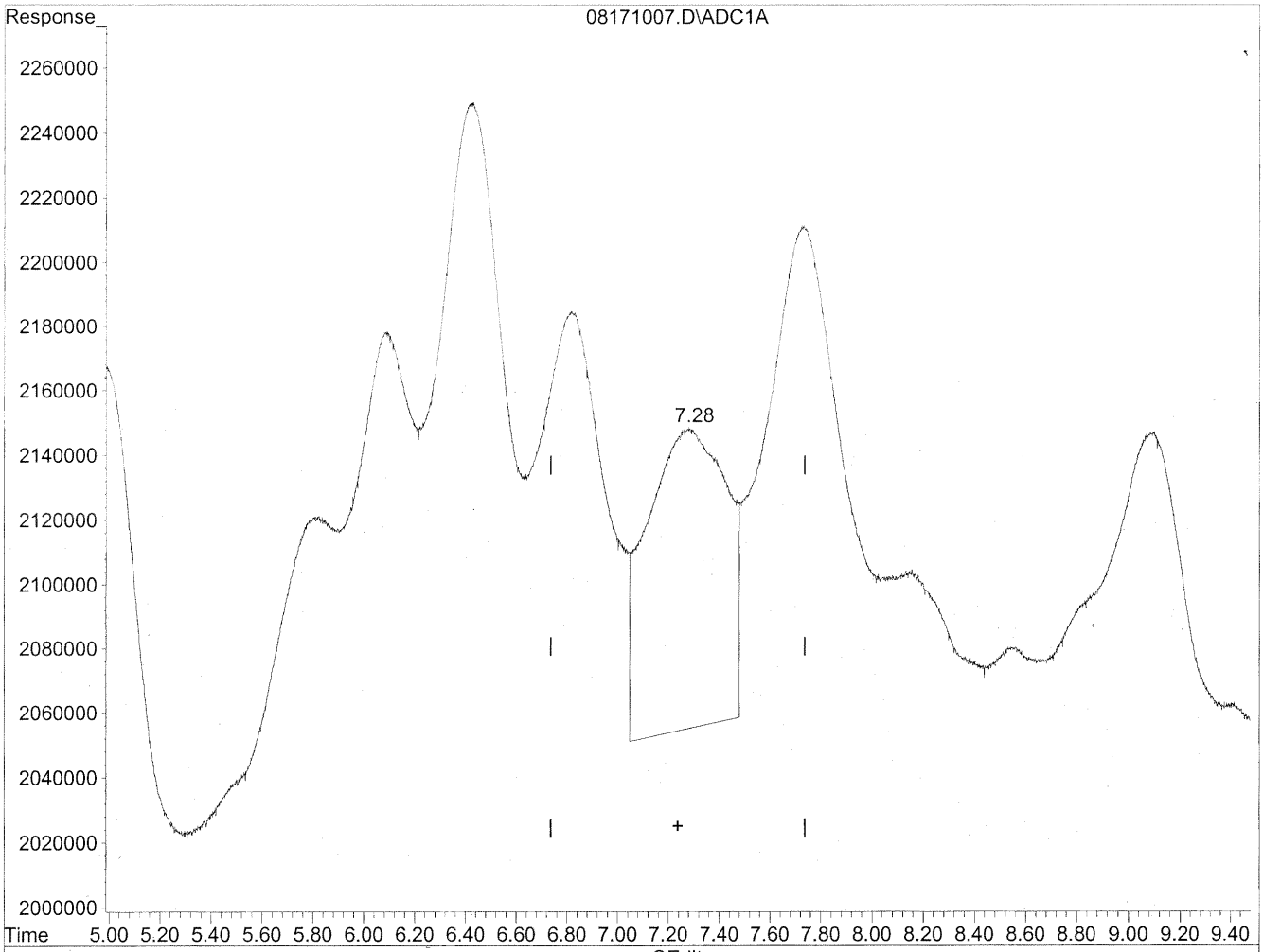
(+) = Expected Retention Time

08171007.D TO110709.M Sat Aug 22 14:10:25 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.28min 259.173ng/ml

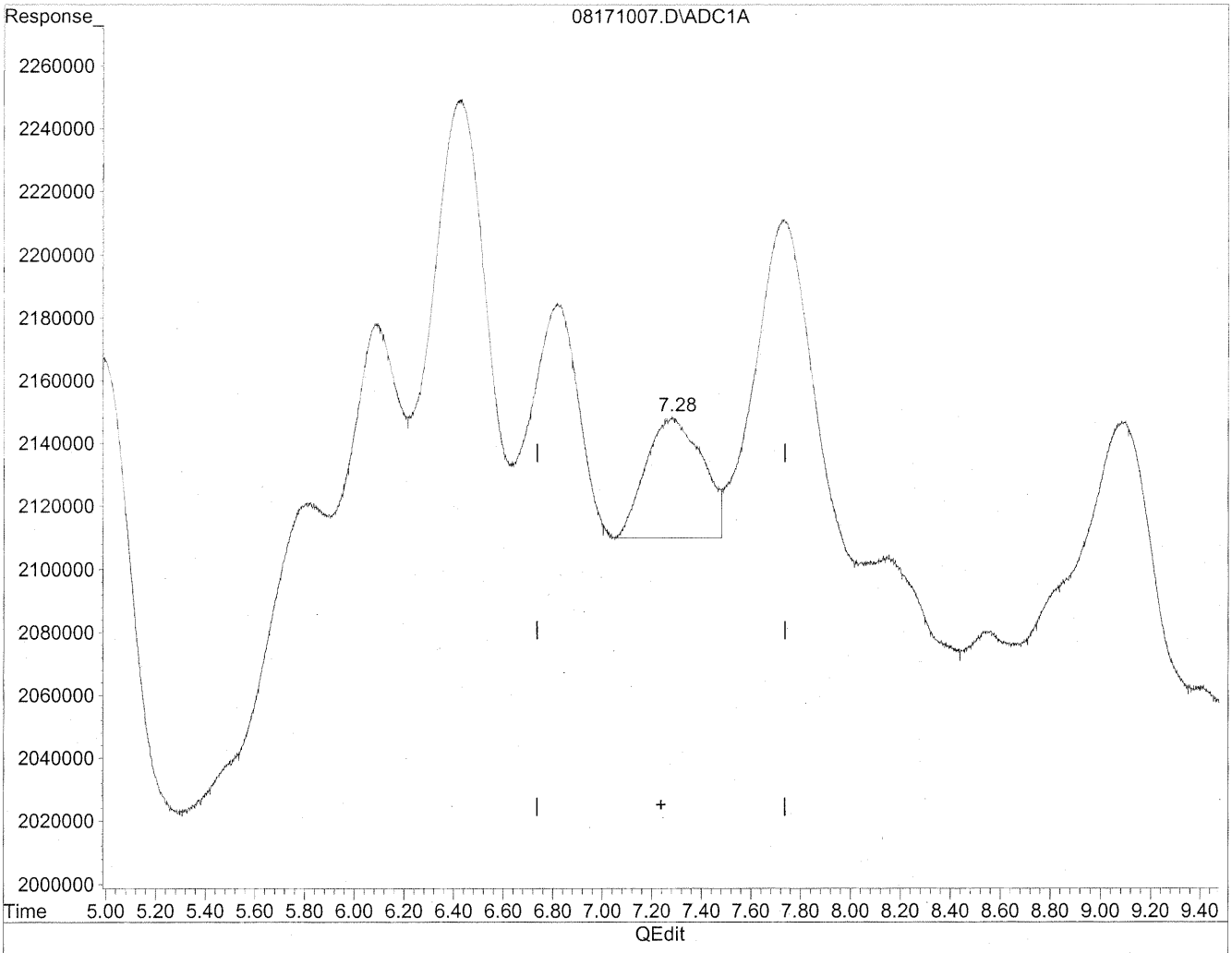
response 20280557

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.28min 77.481ng/ml m

response 6062965

HC 8/22/09
BC
KK 8/23/09

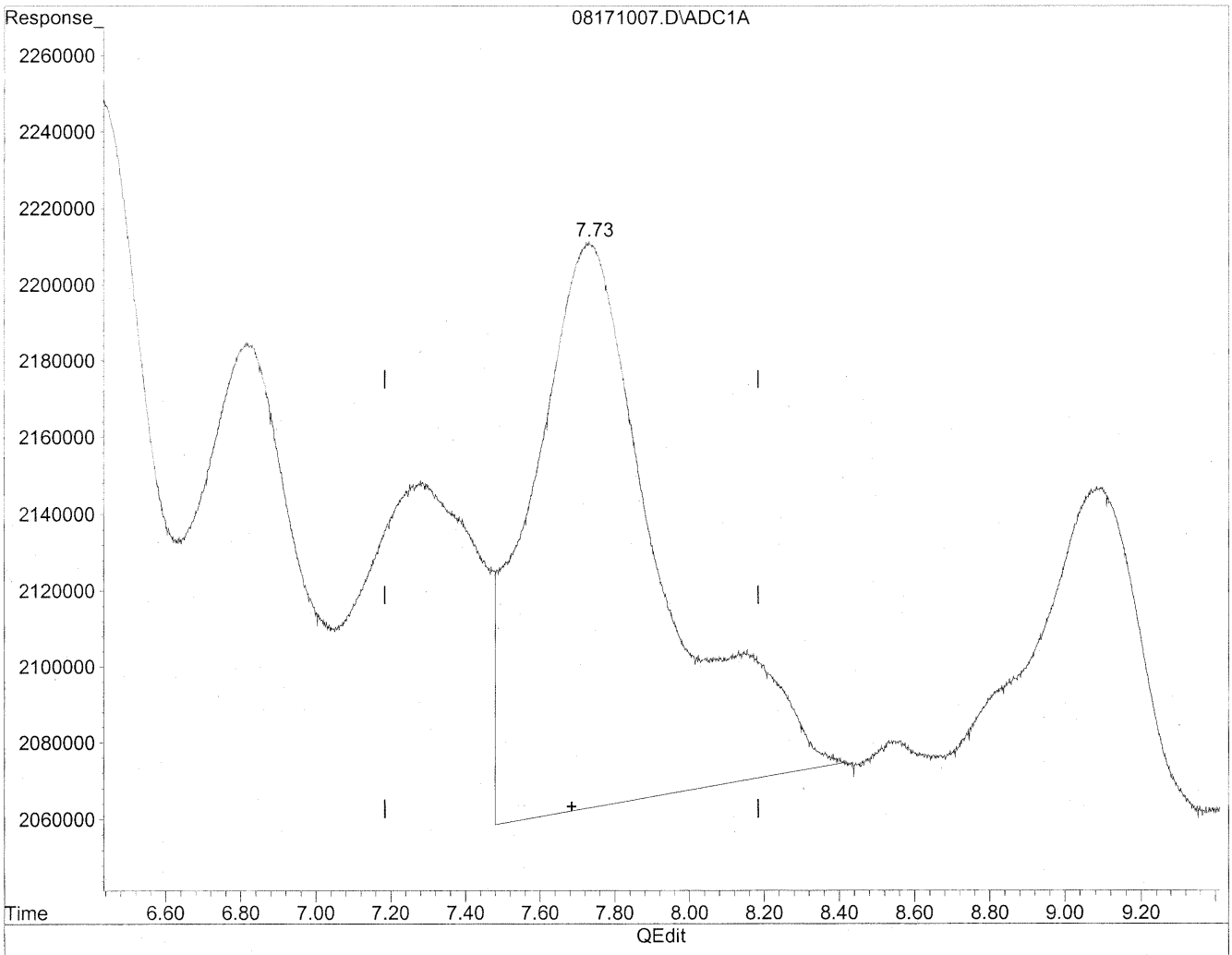
35

(+) = Expected Retention Time
08171007.D TO110709.M Sat Aug 22 14:10:40 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.73min 478.670ng/ml

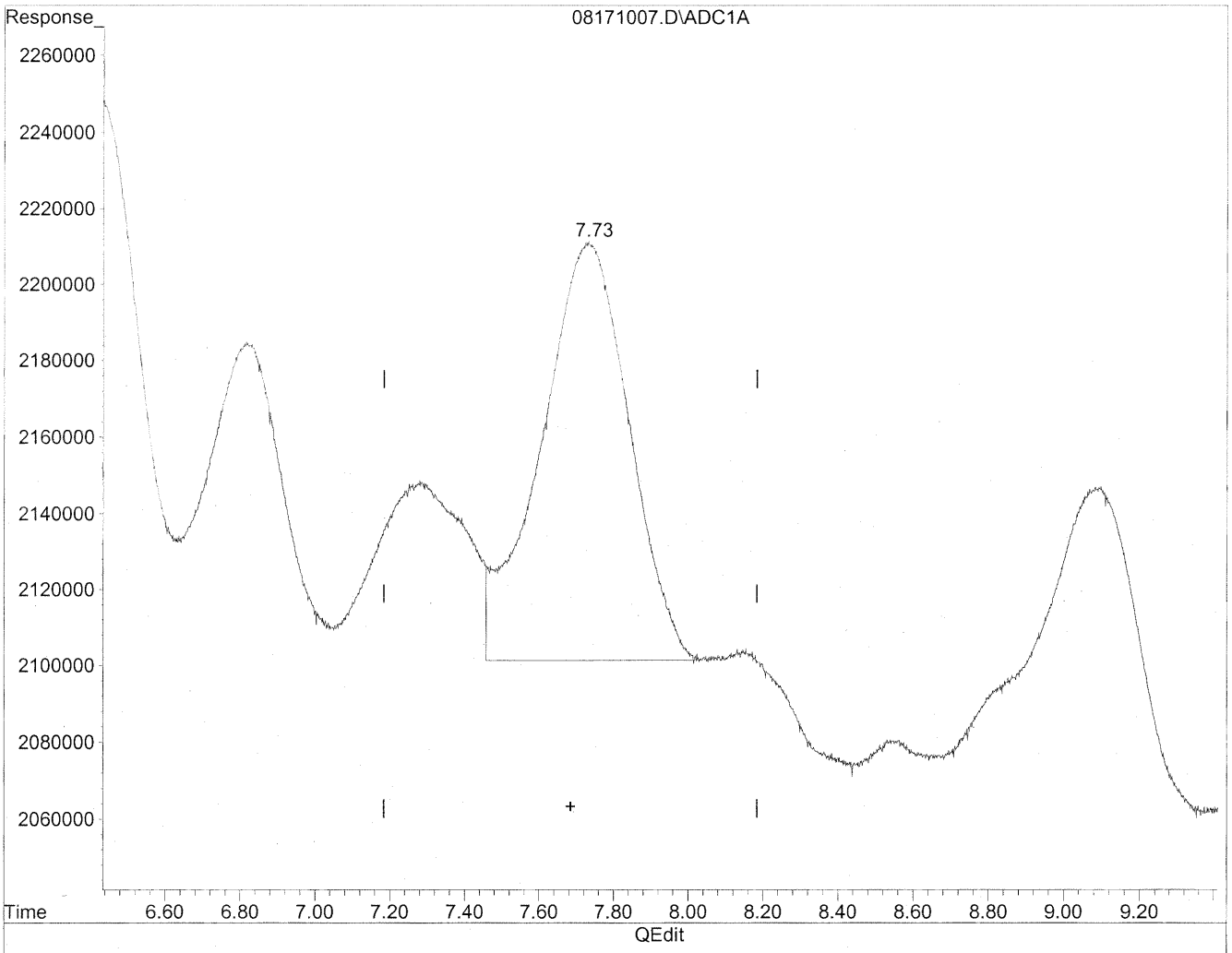
response 35184633

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.73min 245.917ng/ml m

response 18076141

HC 8/22/09

BC 8/23/09

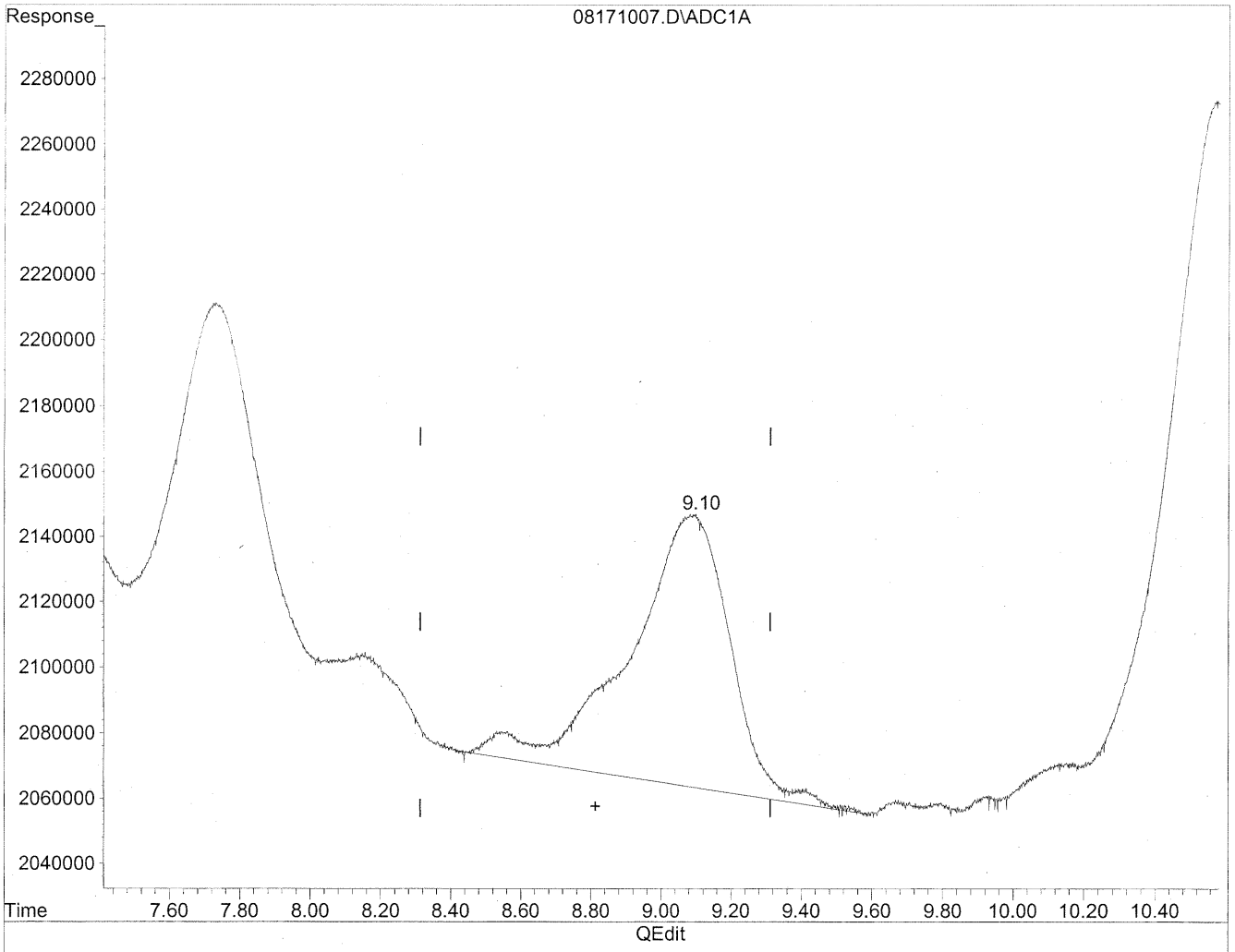
(+) = Expected Retention Time

08171007.D TO110709.M Sat Aug 22 14:11:08 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

9.09min 306.928ng/ml

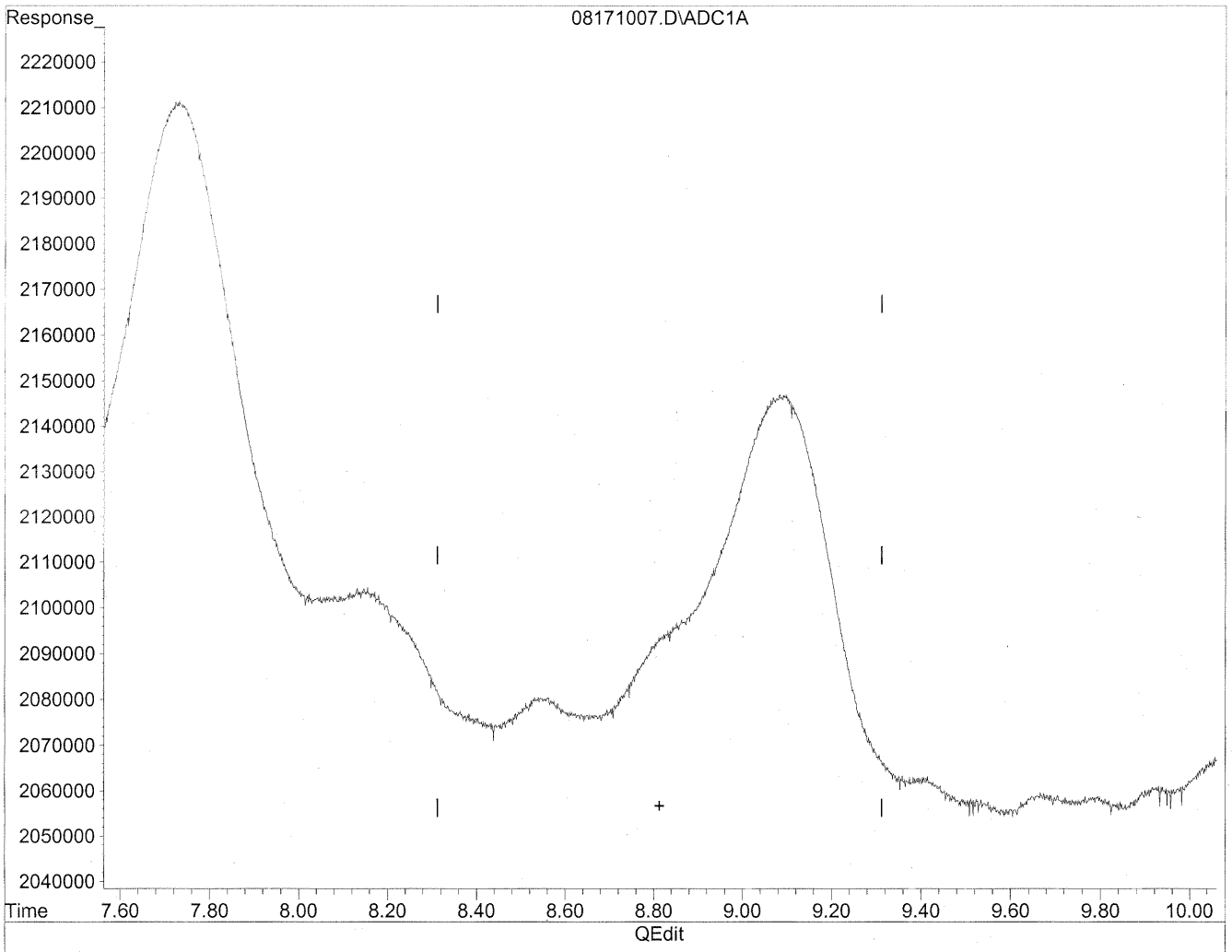
response 16572679

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

0.00min 0.000ng/ml d

response 0

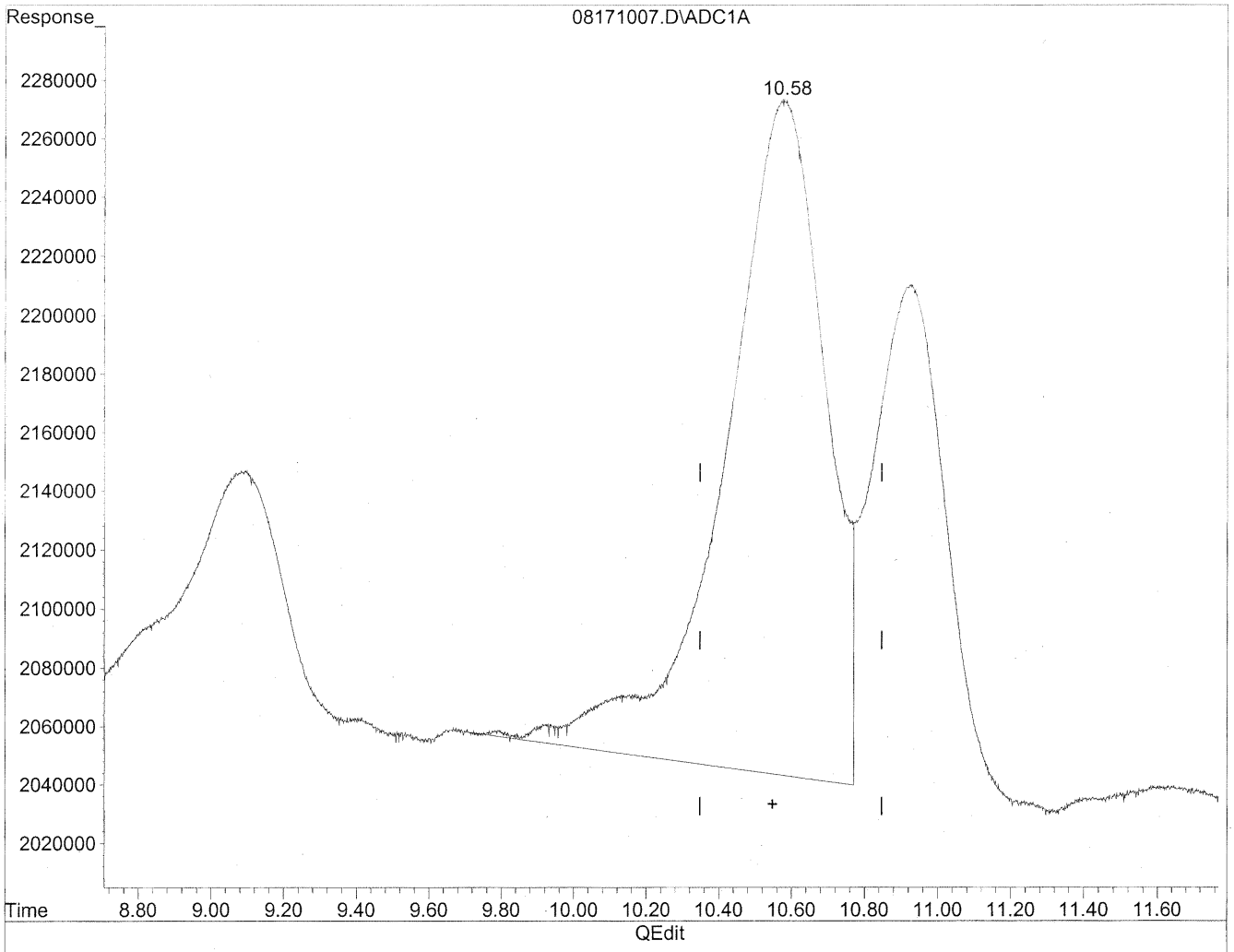
HC 8/22/09
MP
KP 8/23/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.58min 658.652ng/ml

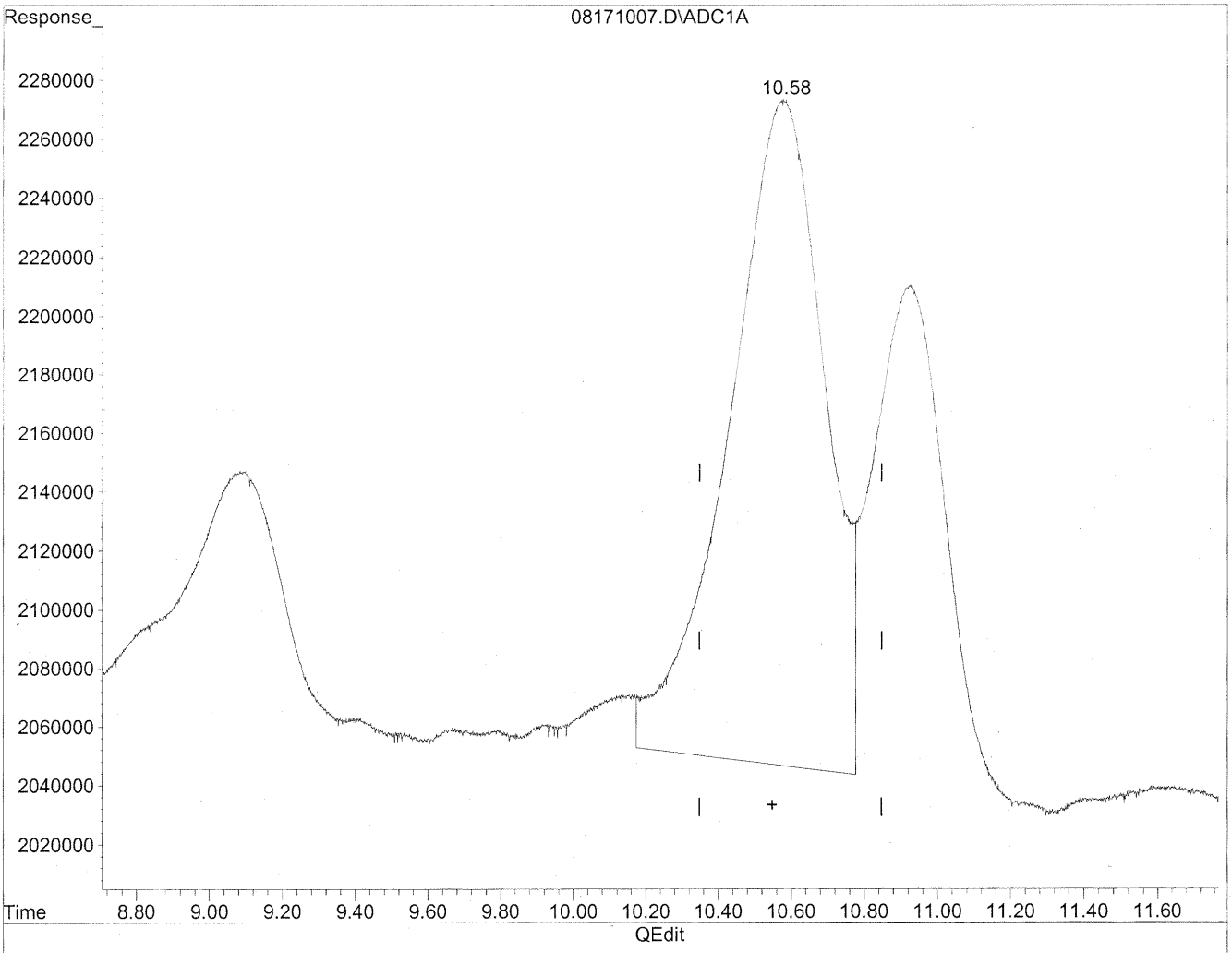
response 44356071

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171007.D Vial: 20
Acq On : 18 Aug 2009 5:24 pm Operator: HC
Sample : P0902772-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:08 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.58min 612.936ng/ml m

response 41277424

hc 8/20/09 SH, BC KP 8/23/09

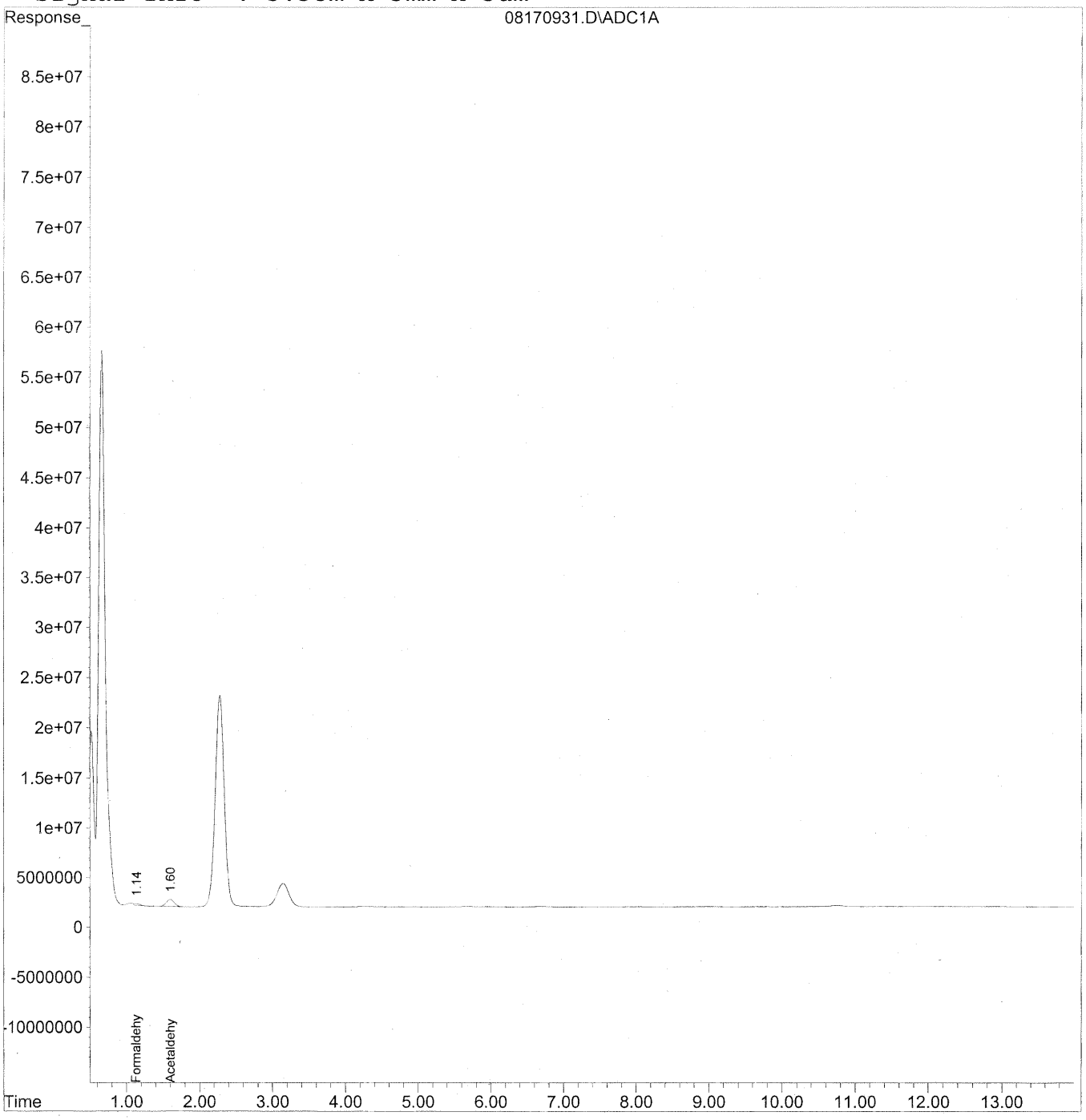
(+) = Expected Retention Time

08171007.D TO110709.M Sat Aug 22 14:11:43 2009

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



42

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

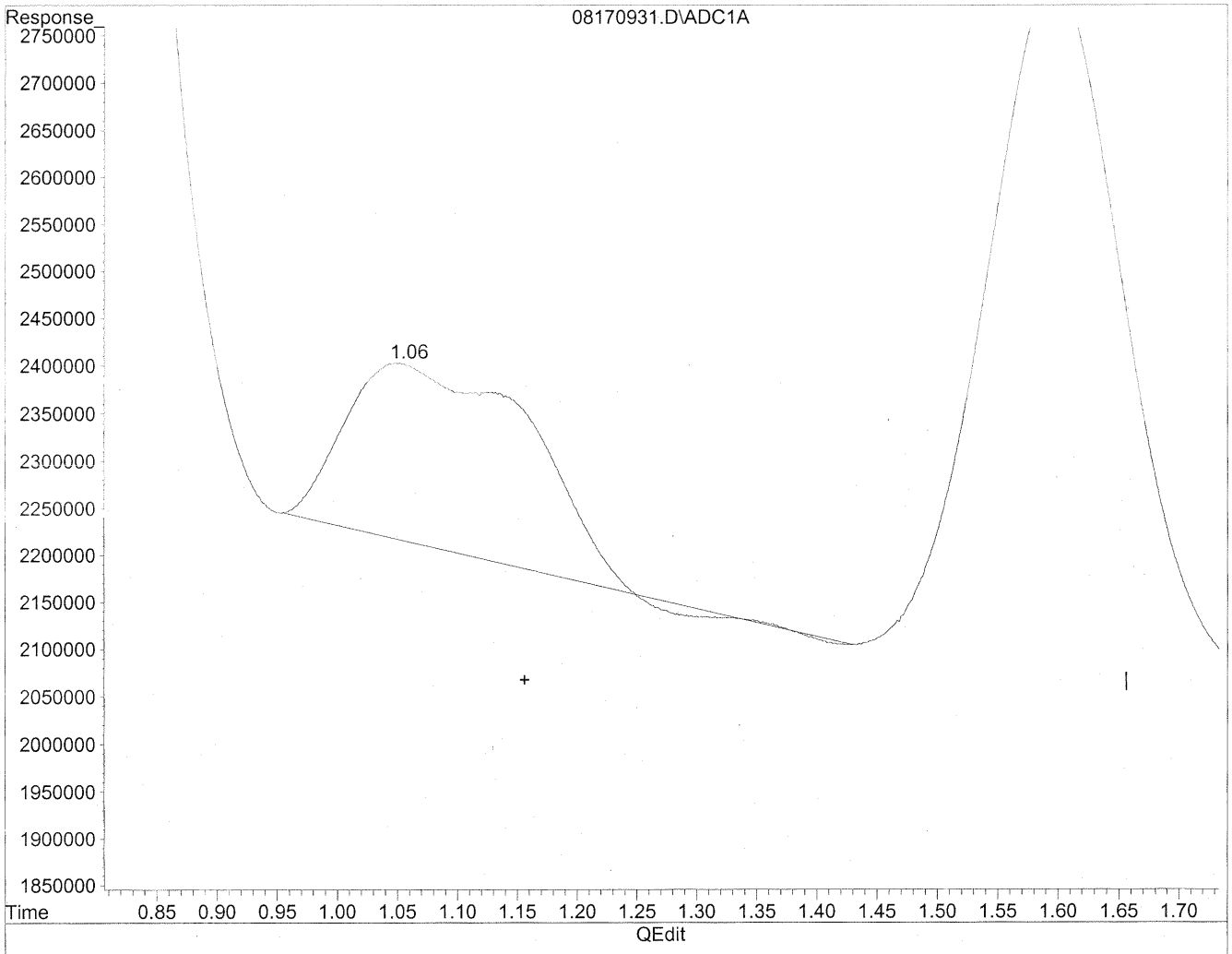
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.14	10879392	59.262	ng/mlm
2)	Acetaldehyde	1.60	56585576	403.538	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/mld
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/mld

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:06 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.05min 112.084ng/ml

response 20576586

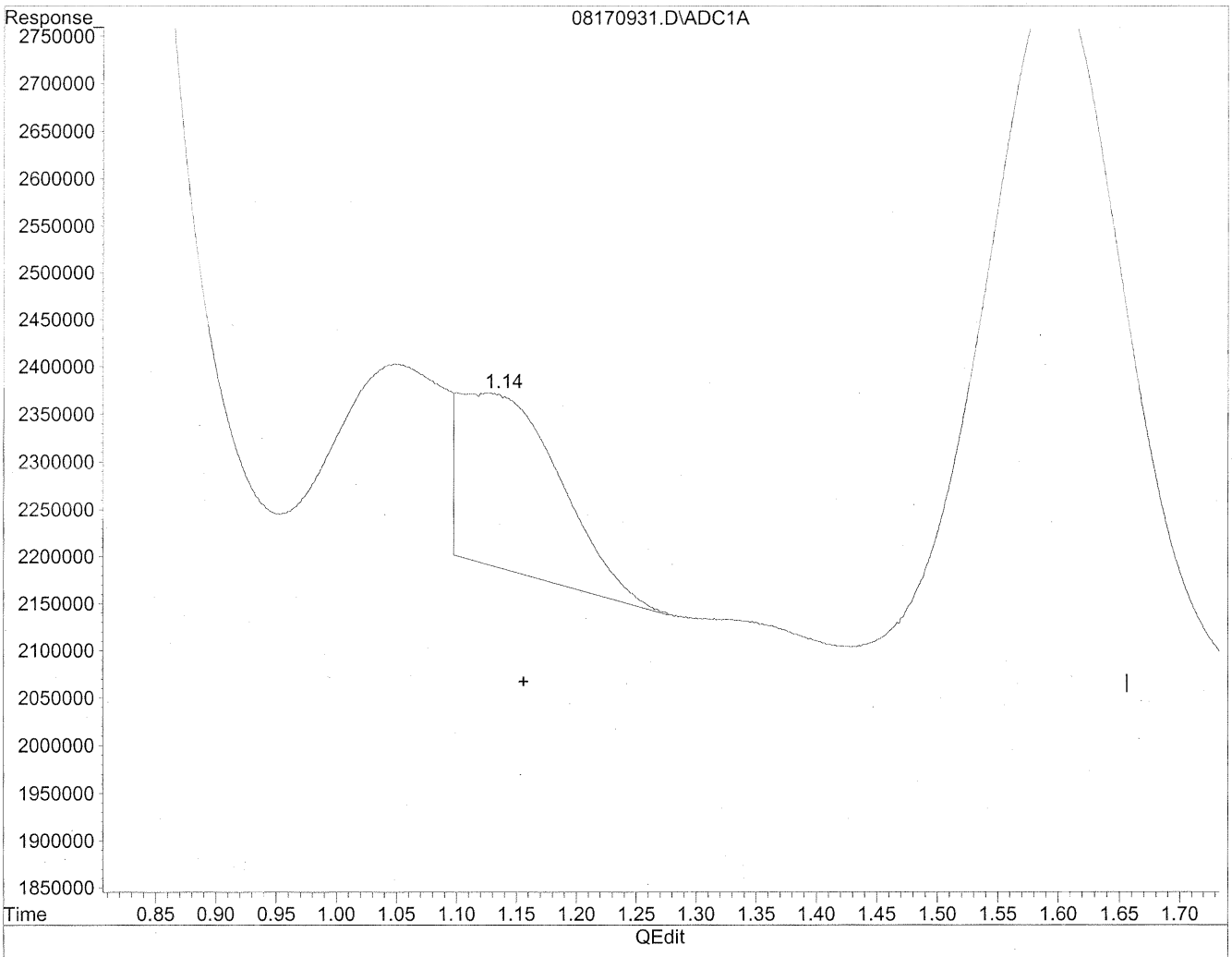
(+) = Expected Retention Time

08170931.D TO110709.M Fri Aug 21 18:06:37 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:06 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.14min 59.262ng/ml m

response 10879392

HC 8/21/09
SP
KK 8/23/09

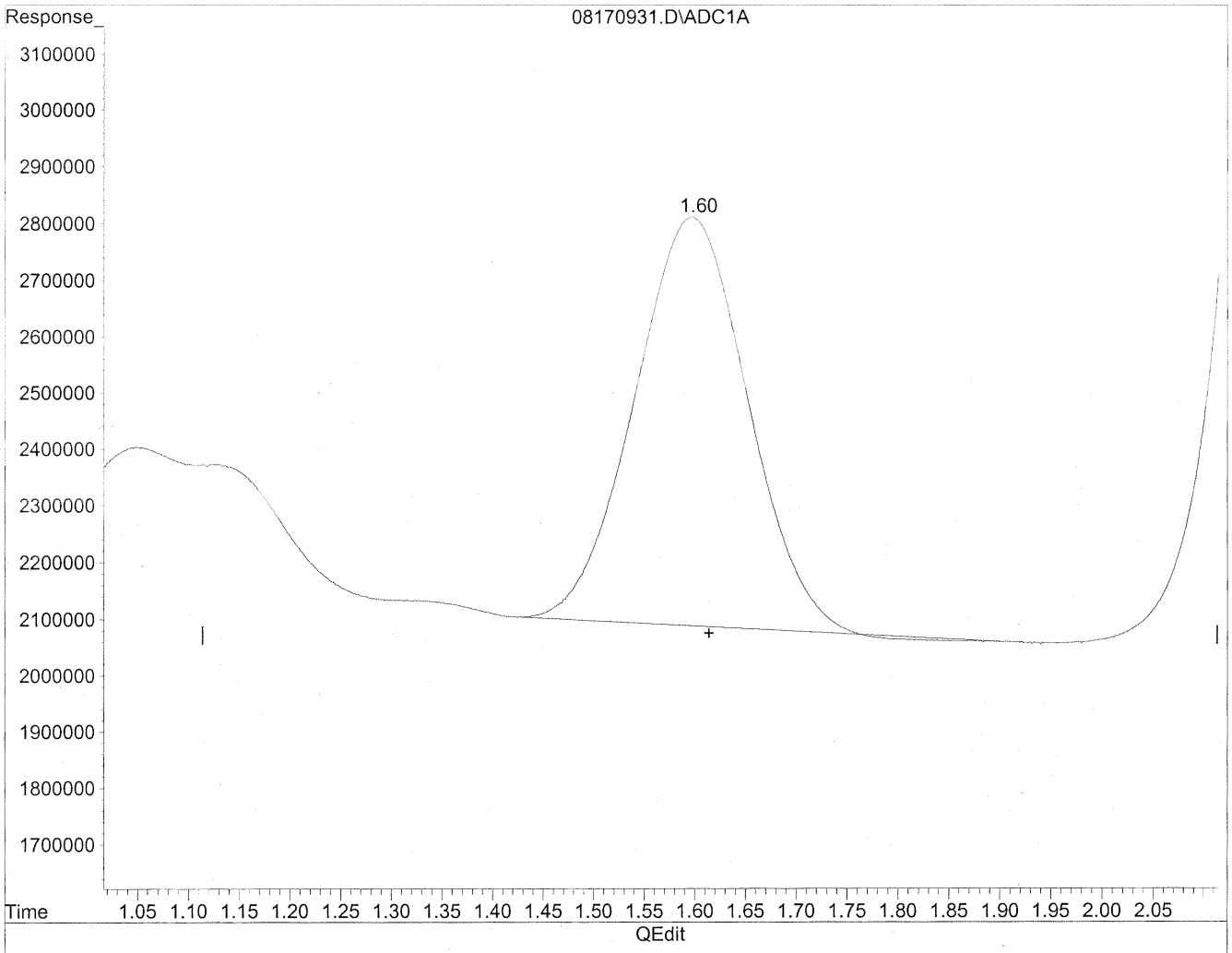
45

(+) = Expected Retention Time
08170931.D TO110709.M Fri Aug 21 18:06:46 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:06 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.60min 398.577ng/ml

response 55889806

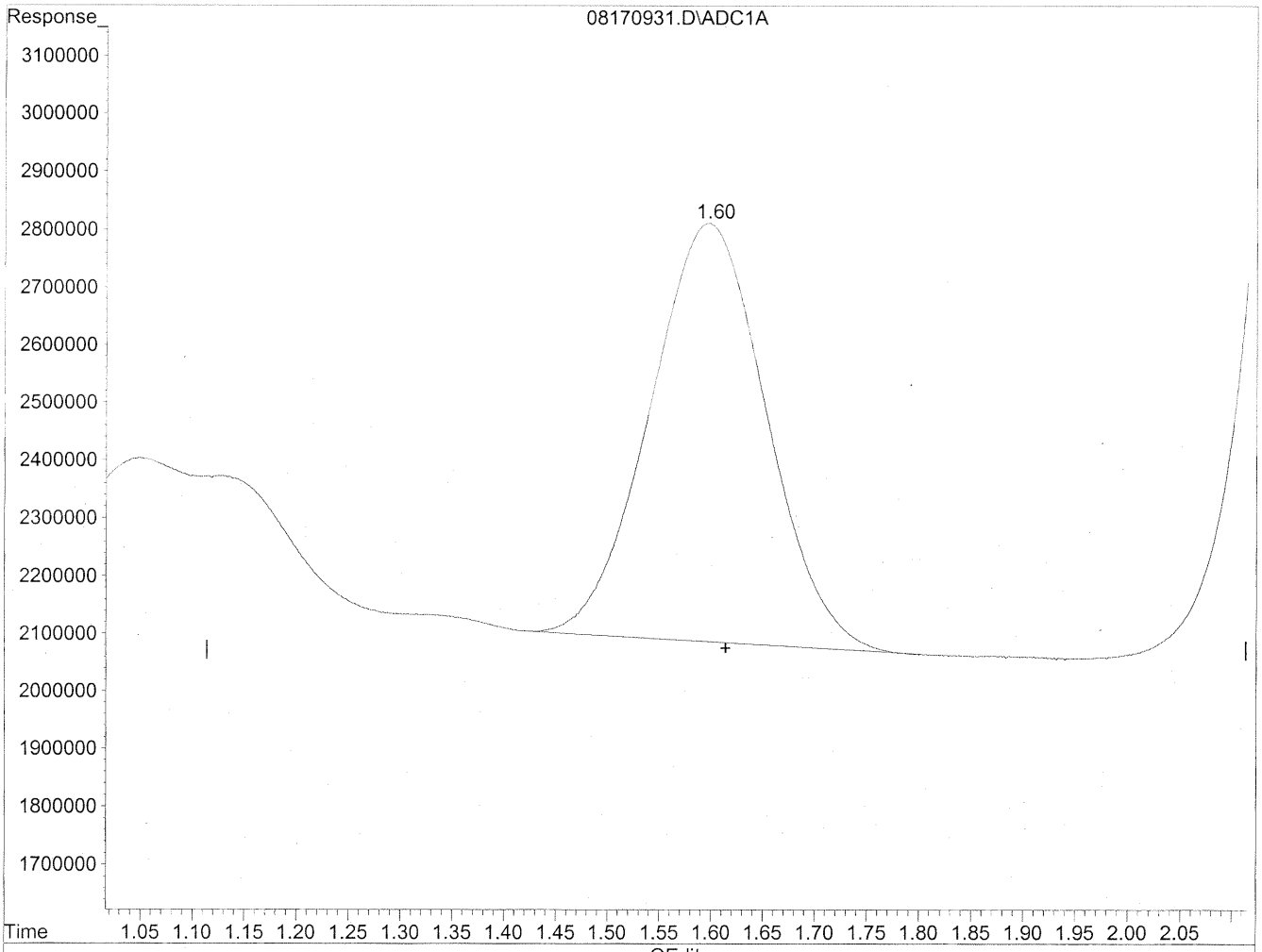
(+) = Expected Retention Time

08170931.D TO110709.M Fri Aug 21 18:06:57 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:06 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(2) Acetaldehyde

1.60min 403.538ng/ml m

response 56585576

HC 8/21/09
MP
KK 8/23/09

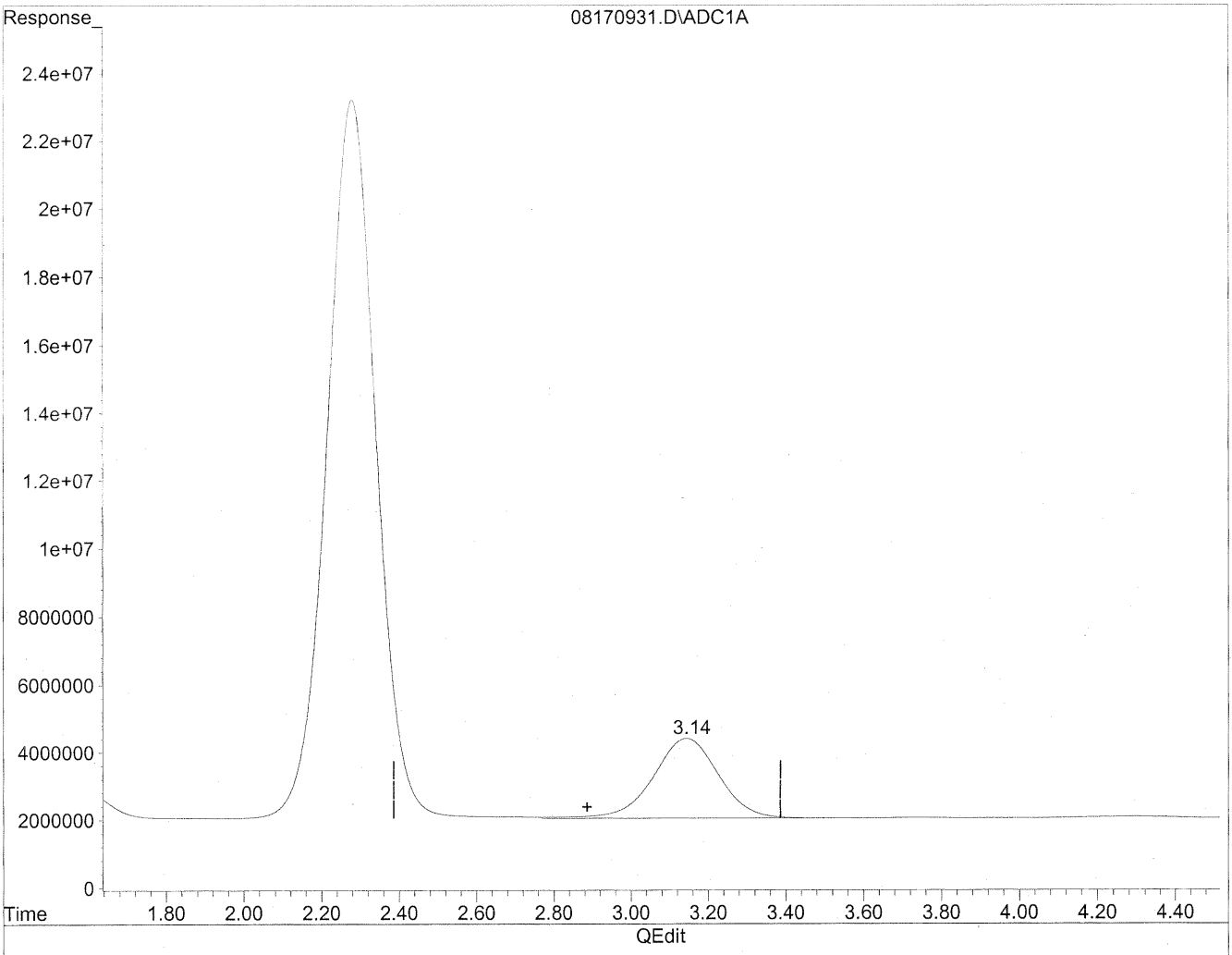
47

(+) = Expected Retention Time
08170931.D TO110709.M Fri Aug 21 18:07:03 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:06 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.14min 2602.477ng/ml

response 277671857

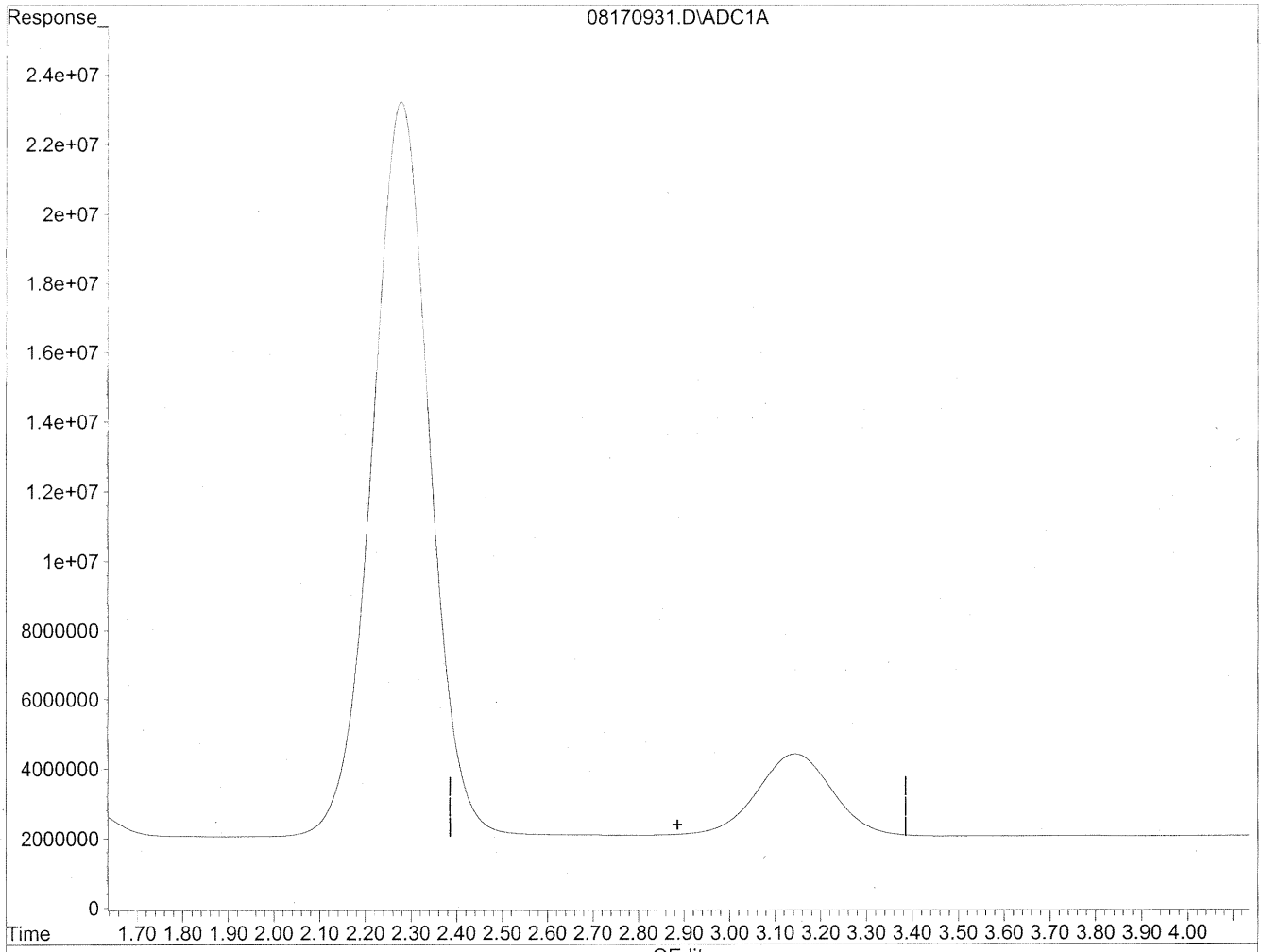
(+) = Expected Retention Time

08170931.D TO110709.M Fri Aug 21 18:07:13 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:06 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/21/09 MP

KK 8/23/09

49

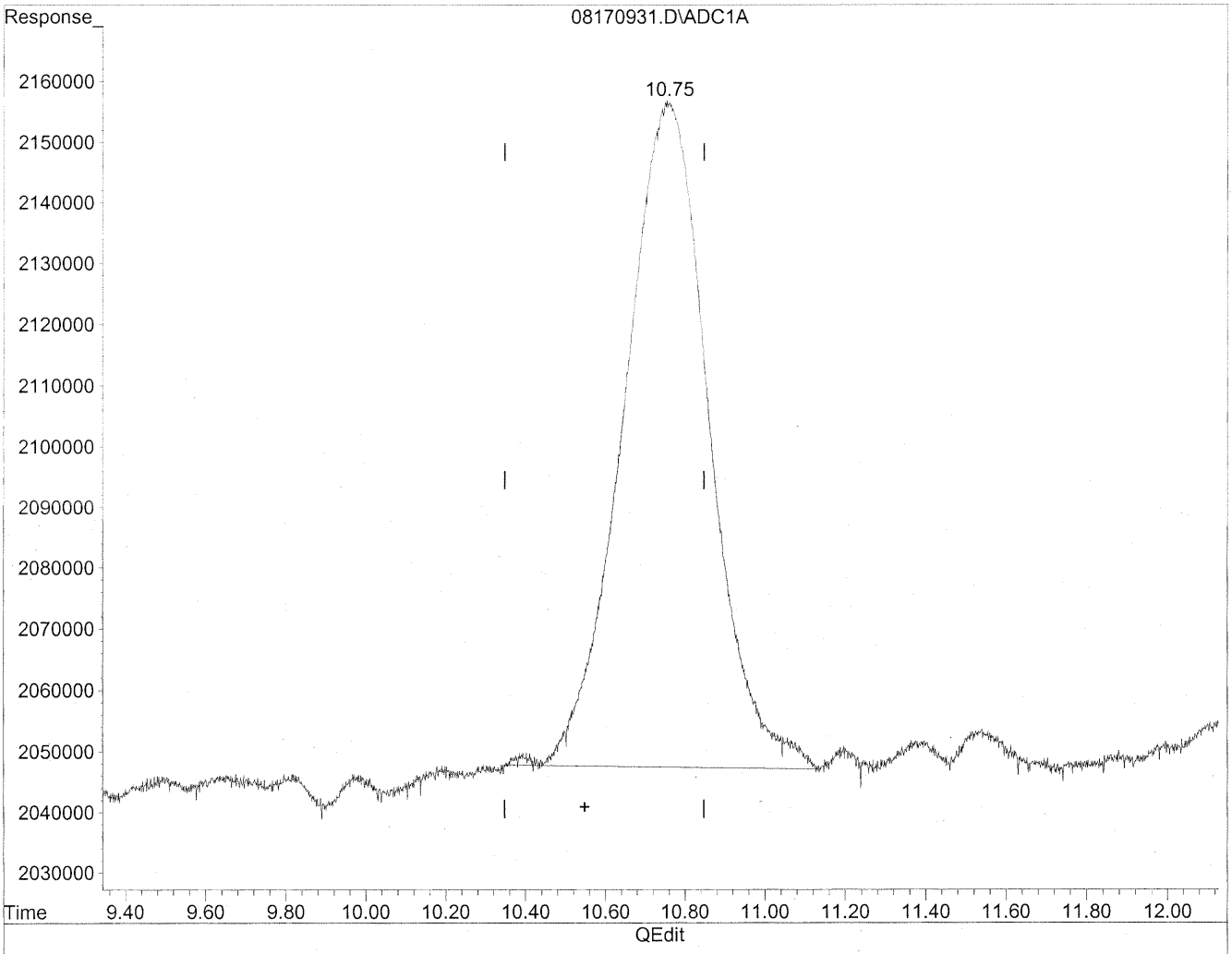
(+) = Expected Retention Time

08170931.D TO110709.M Fri Aug 21 18:07:40 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:06 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.76min 240.900ng/ml

response 16223144

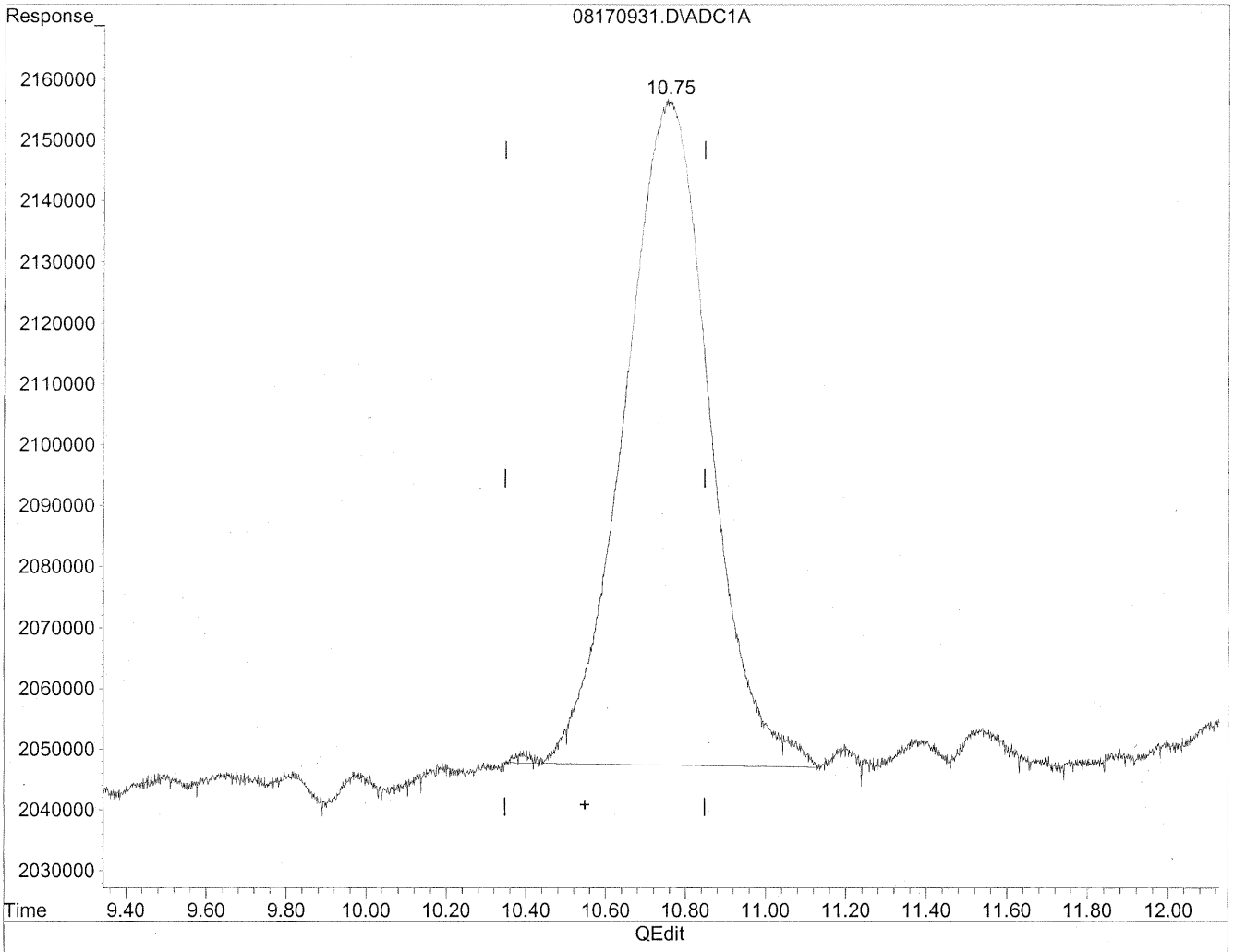
(+) = Expected Retention Time

08170931.D TO110709.M Fri Aug 21 18:07:46 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:06 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.76min 240.900ng/ml

response 16223144

HC 8/21/09 TC

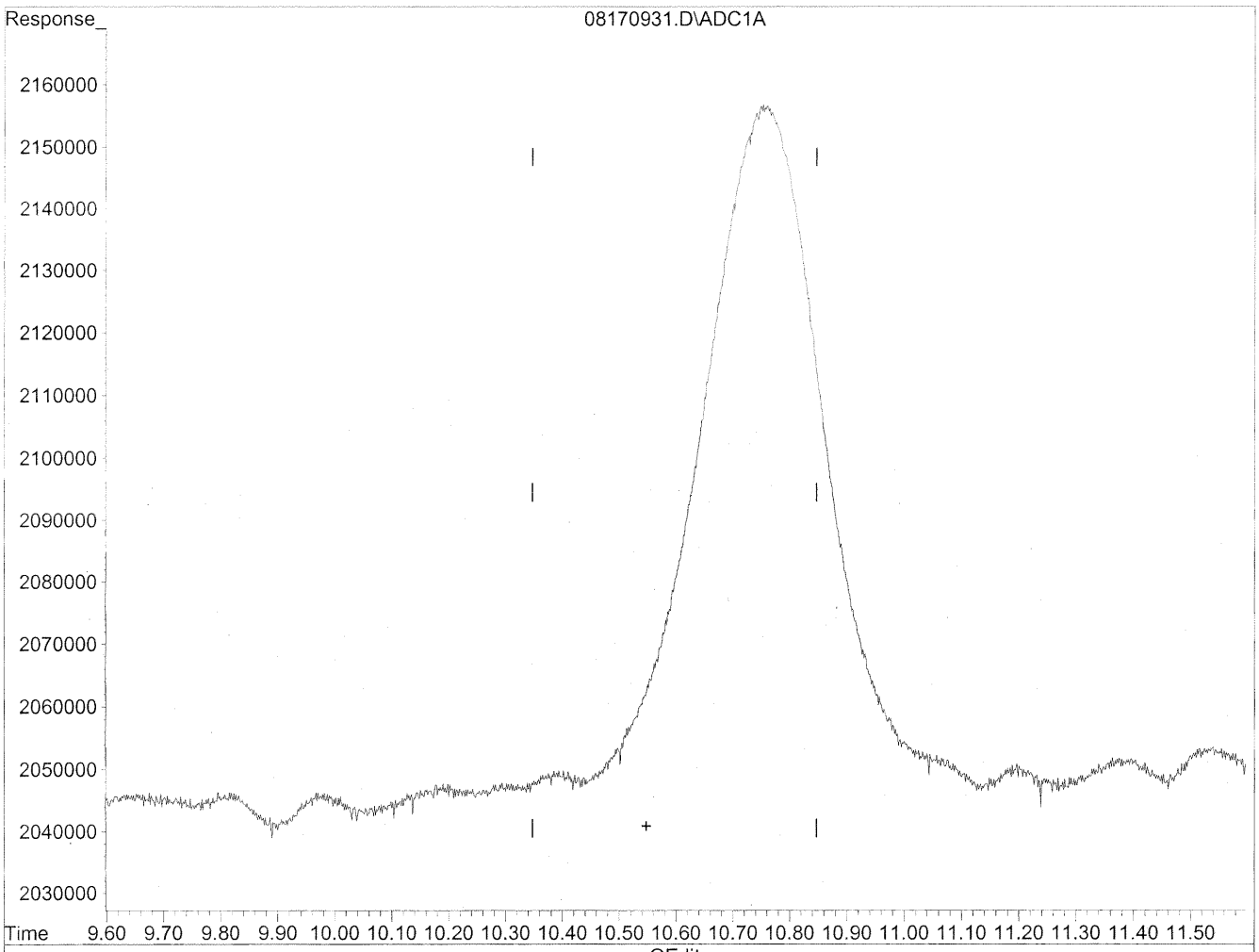
(+) = Expected Retention Time

08170931.D TO110709.M Fri Aug 21 18:12:04 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/25/09

MP

KP 8/25/09

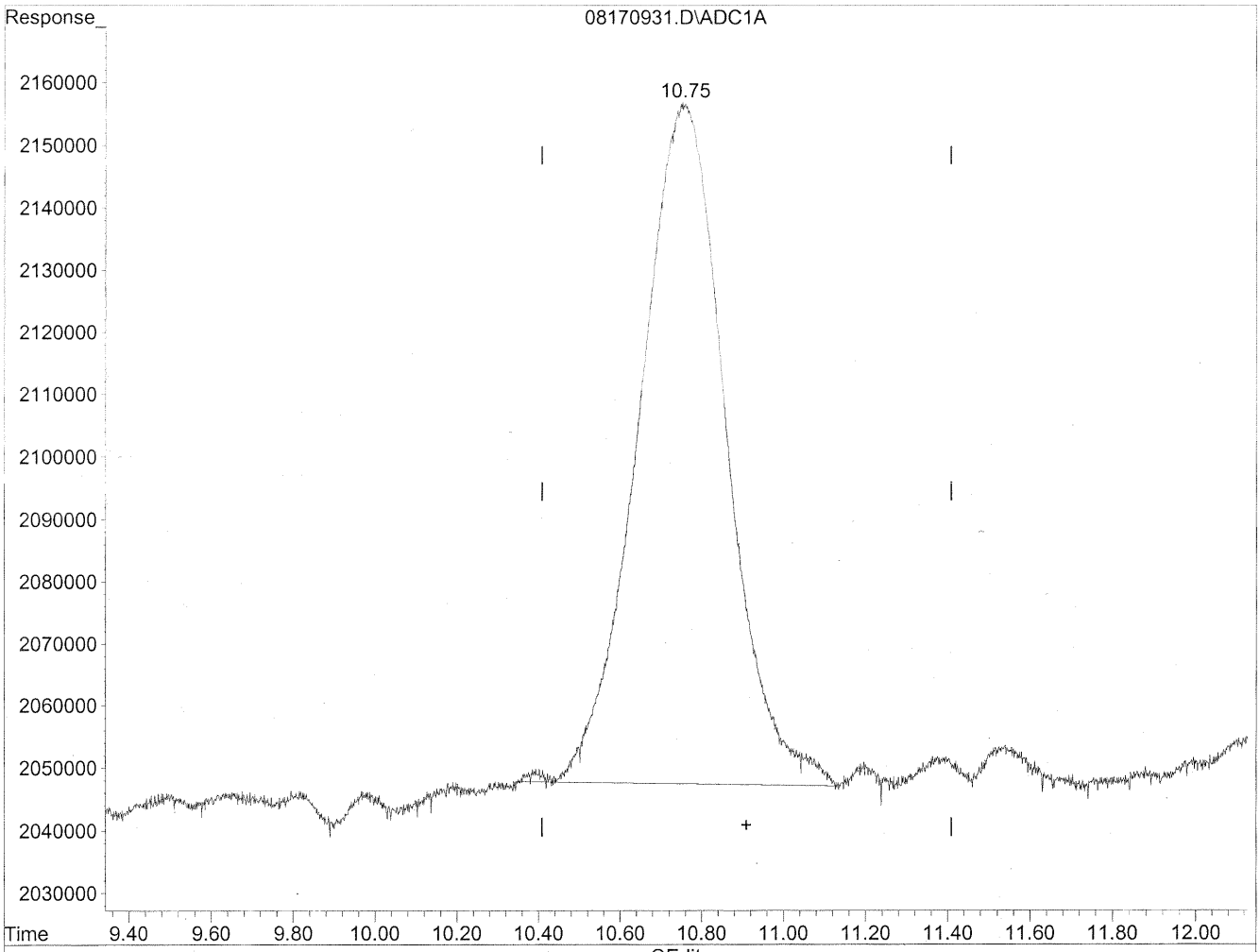
52

(+) = Expected Retention Time
08170931.D TO110709.M Tue Aug 25 07:39:24 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:12 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

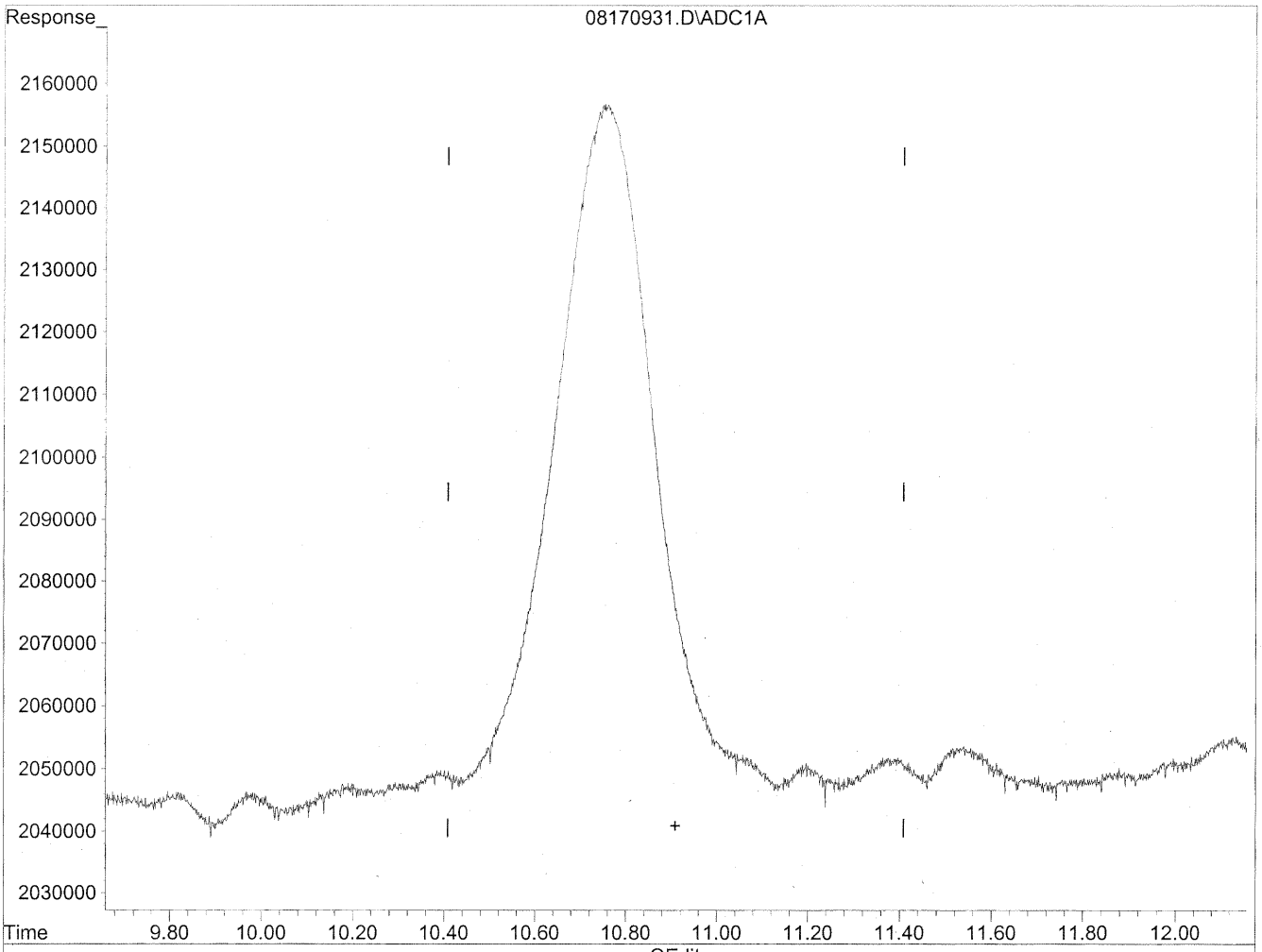
10.76min 330.994ng/ml

response 16223144

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:12 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/21/09

MP

KR 8/23/09

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **100305**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P0902772-003

Test Code:	EPA Method TO-11A	Date Collected:	8/10/09
Instrument ID:	Waters LC Module I Plus/UV_Vis 360/LC1	Date Received:	8/12/09
Analyst:	Hani Cherazaie	Date Analyzed:	8/17 - 8/18/09
Sampling Media:	Silica Gel DNPH Tube	Desorption Volume:	1.0 ml
Test Notes:	**BC**	Volume Sampled:	106 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**3,700**	**35**	0.94	**29**	0.77	
75-07-0	**Acetaldehyde**	**2,600**	**24**	0.94	**13**	0.52	**BT**
123-38-6	**Propionaldehyde**	**220**	**2.1**	0.94	**0.87**	0.40	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.94	ND	0.33	
123-72-8	**Butyraldehyde**	**220**	**2.1**	0.94	**0.71**	0.32	
100-52-7	**Benzaldehyde**	**340**	**3.3**	0.94	**0.75**	0.22	
590-86-3	Isovaleraldehyde	< 100	ND	0.94	ND	0.27	
110-62-3	**Valeraldehyde**	**240**	**2.2**	0.94	**0.64**	0.27	
529-20-4	o-Tolualdehyde	< 100	ND	0.94	ND	0.19	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	1.9	ND	0.38	
66-25-1	**n-Hexaldehyde**	**780**	**7.4**	0.94	**1.8**	0.23	
5779-94-2	**2,5-Dimethylbenzaldehyde**	**740**	**7.0**	0.94	**1.3**	0.17	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
BT = Results indicated possible breakthrough; back section > 10% front section.

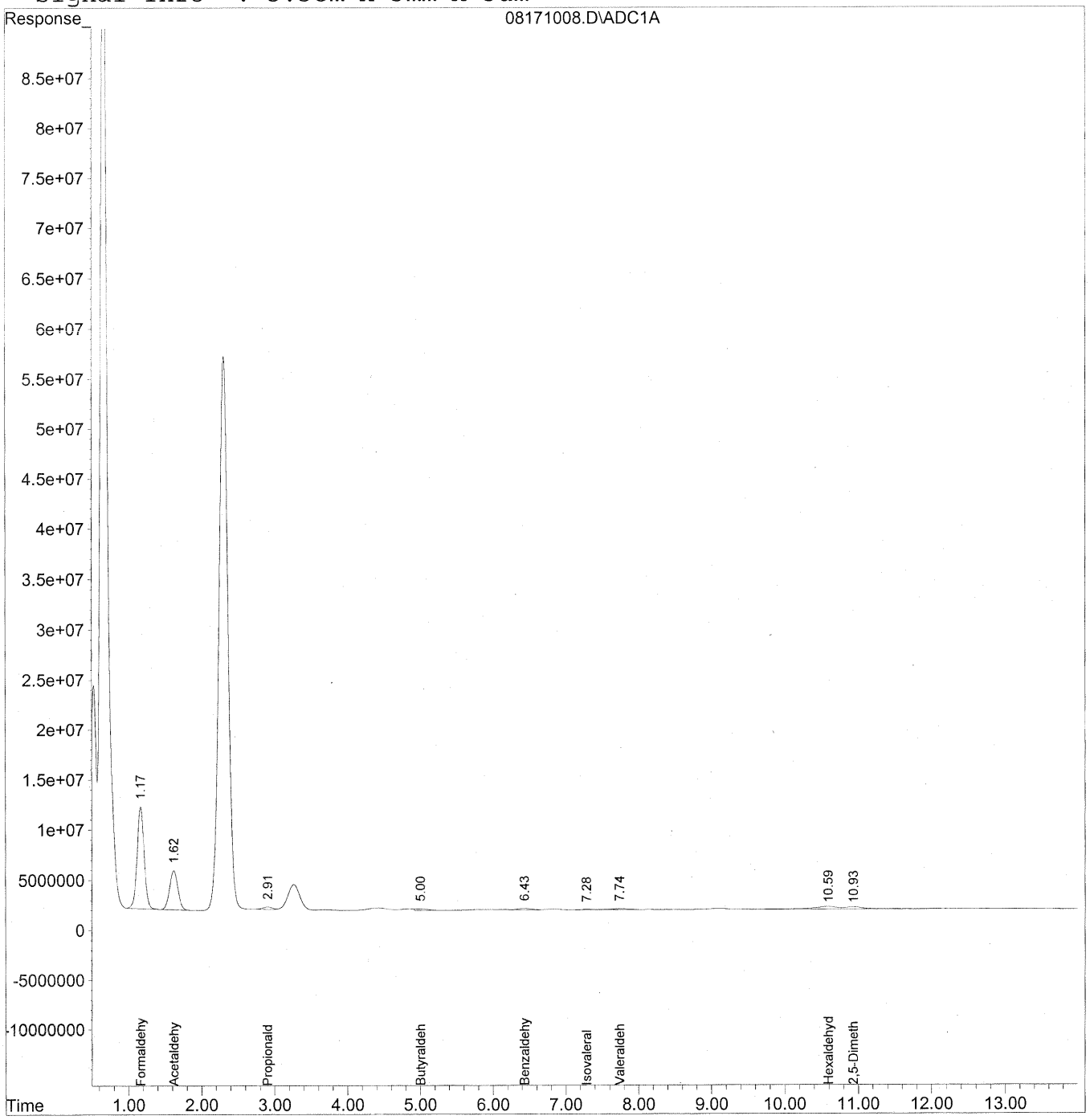
Verified By: _____ Date: 8/25/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:15 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



56

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:15 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

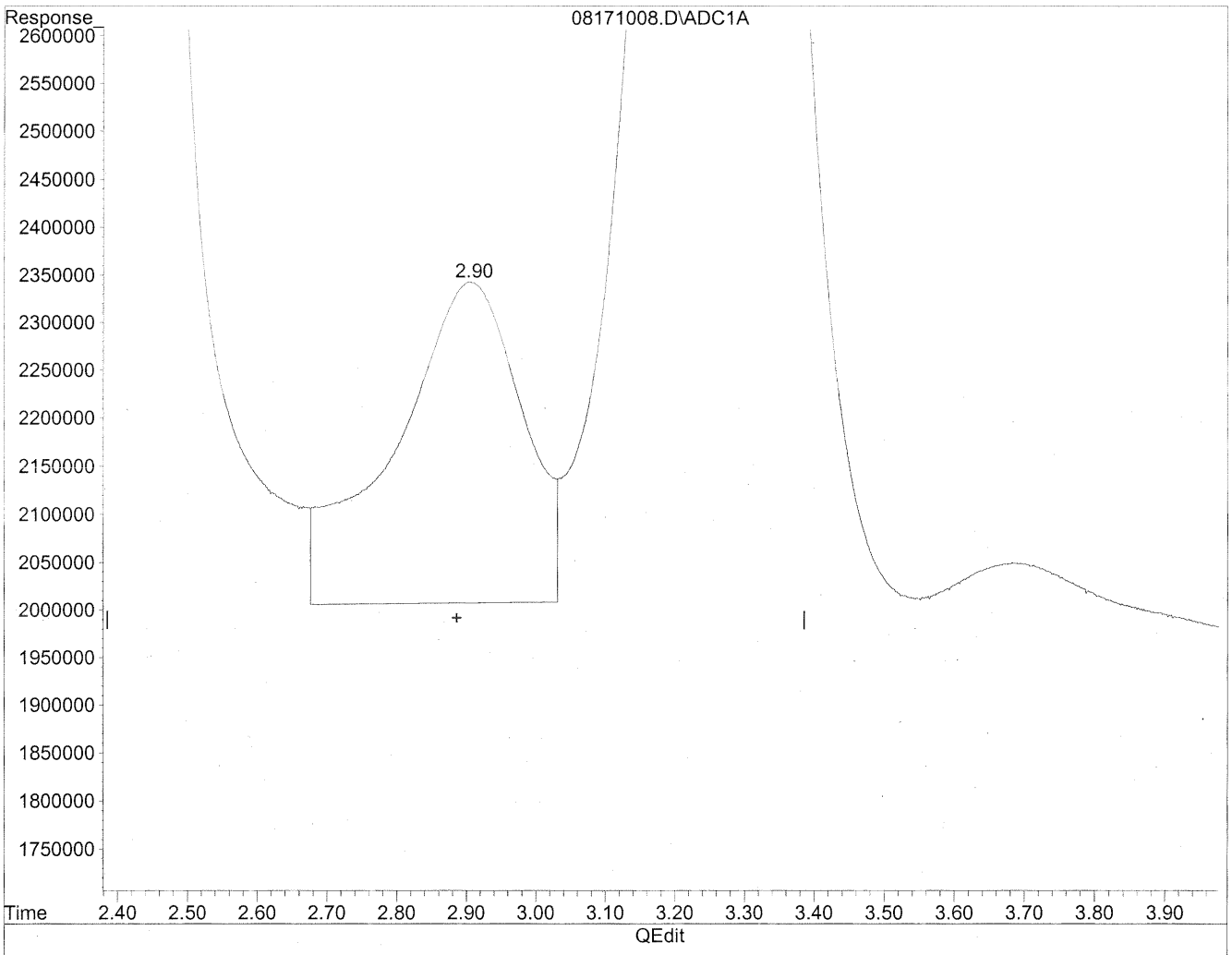
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	682988313	3720.357	ng/ml
2)	Acetaldehyde	1.62	311745615	2223.205	ng/ml
3)	Propionaldehyde	2.91	23430095	219.598	ng/mlm
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	5.00	19731484	223.368	ng/mlm
6)	Benzaldehyde	6.43	22696600	344.570	ng/mlm
7)	Isovaleraldehyde	7.28	6362152	81.304	ng/mlm
8)	Valeraldehyde	7.74	17523823	238.403	ng/mlm
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/mld
11)	Hexaldehyde	10.59	52627599	781.477	ng/mlm
12)	2,5-Dimethylbenzaldehyde	10.93	36497004	744.633	ng/mlm

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.91min 399.002ng/ml

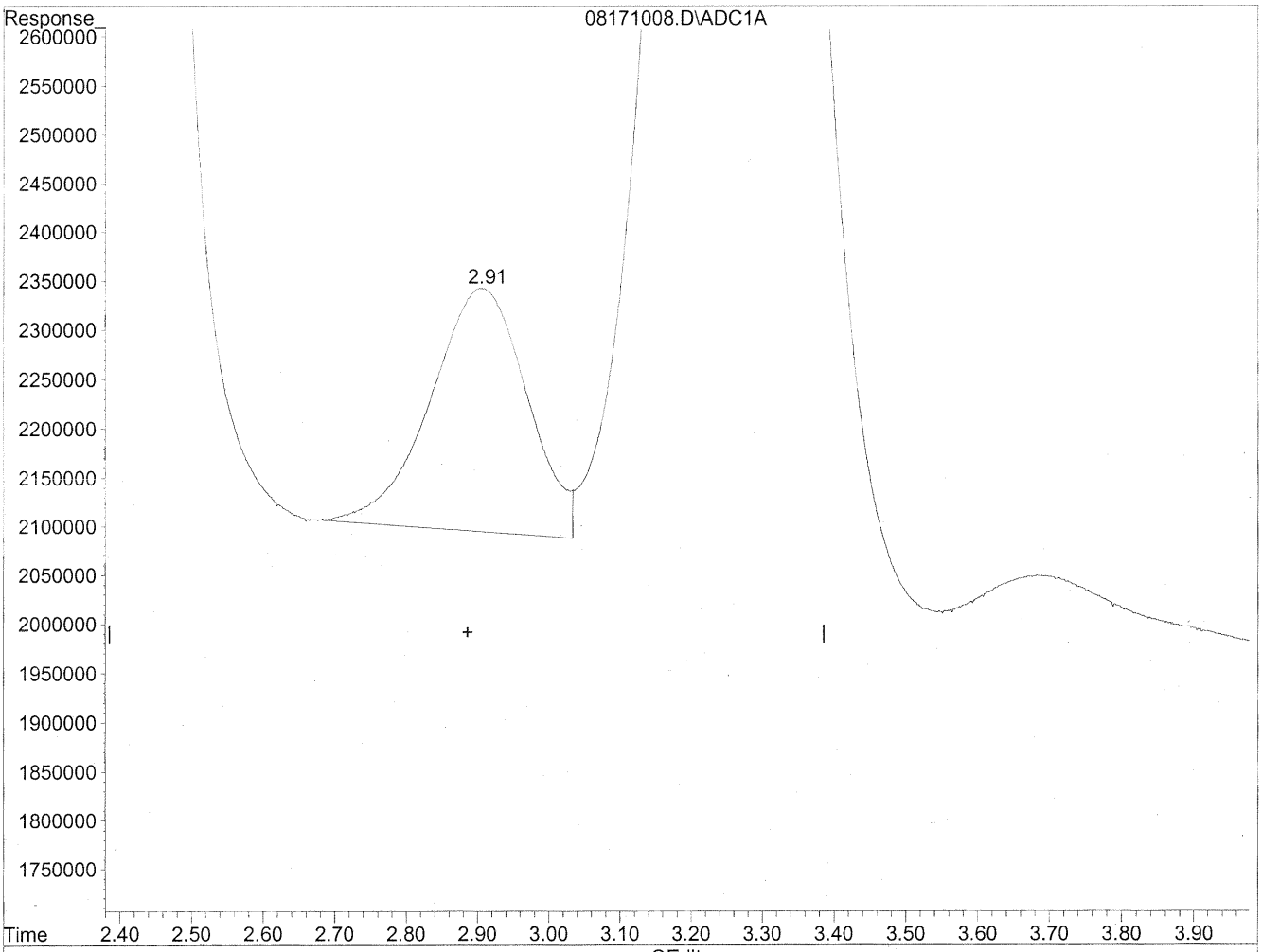
response 42571603

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.91min 219.598ng/ml m

response 23430095

HC 8/22/09
BL
KR 8/23/09

59

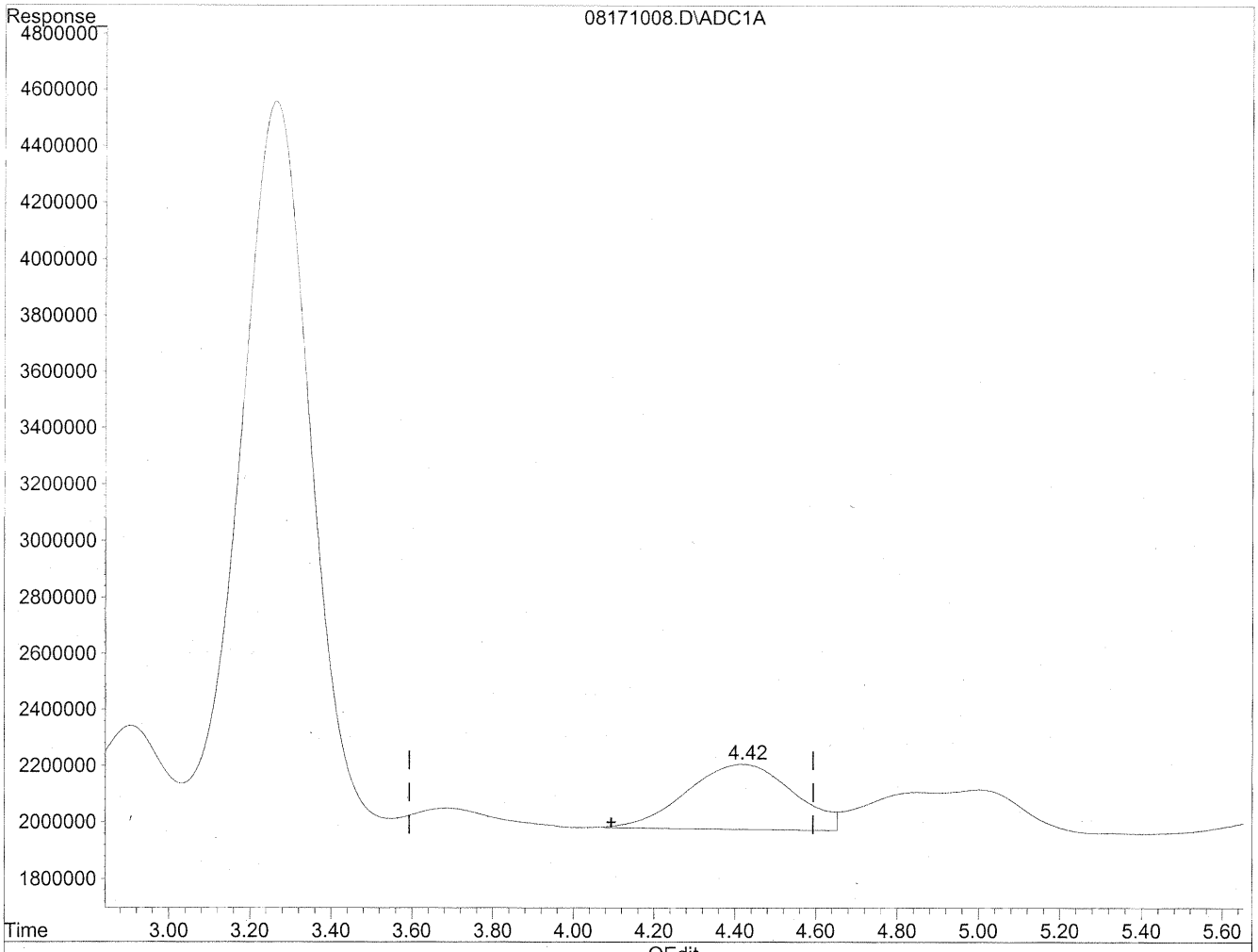
(+) = Expected Retention Time

08171008.D TO110709.M Sat Aug 22 14:13:06 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.42min 434.908ng/ml

response 42366684

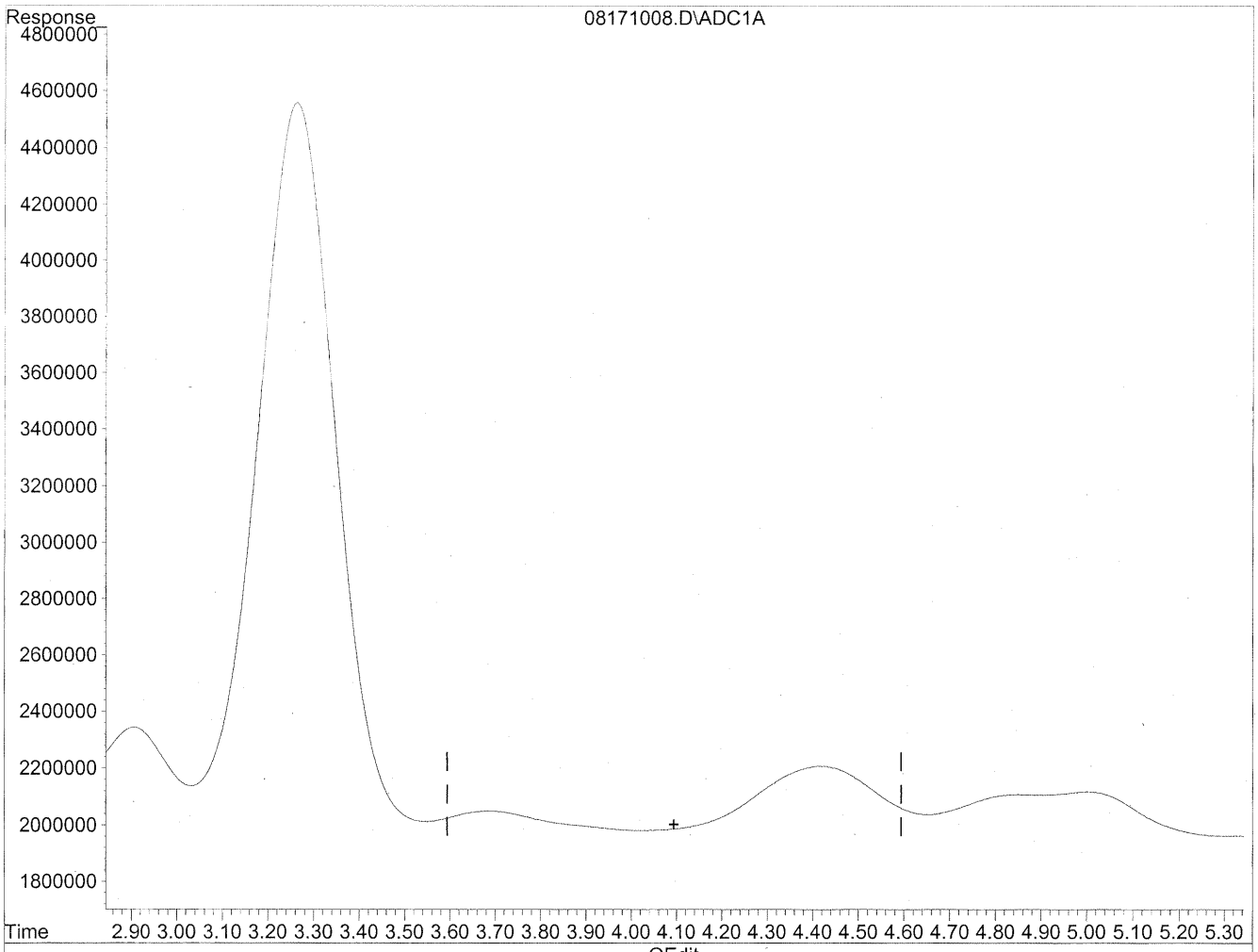
(+) = Expected Retention Time

08171008.D TO110709.M Sat Aug 22 14:13:10 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC
8/22/09
MP
KR 8/23/09

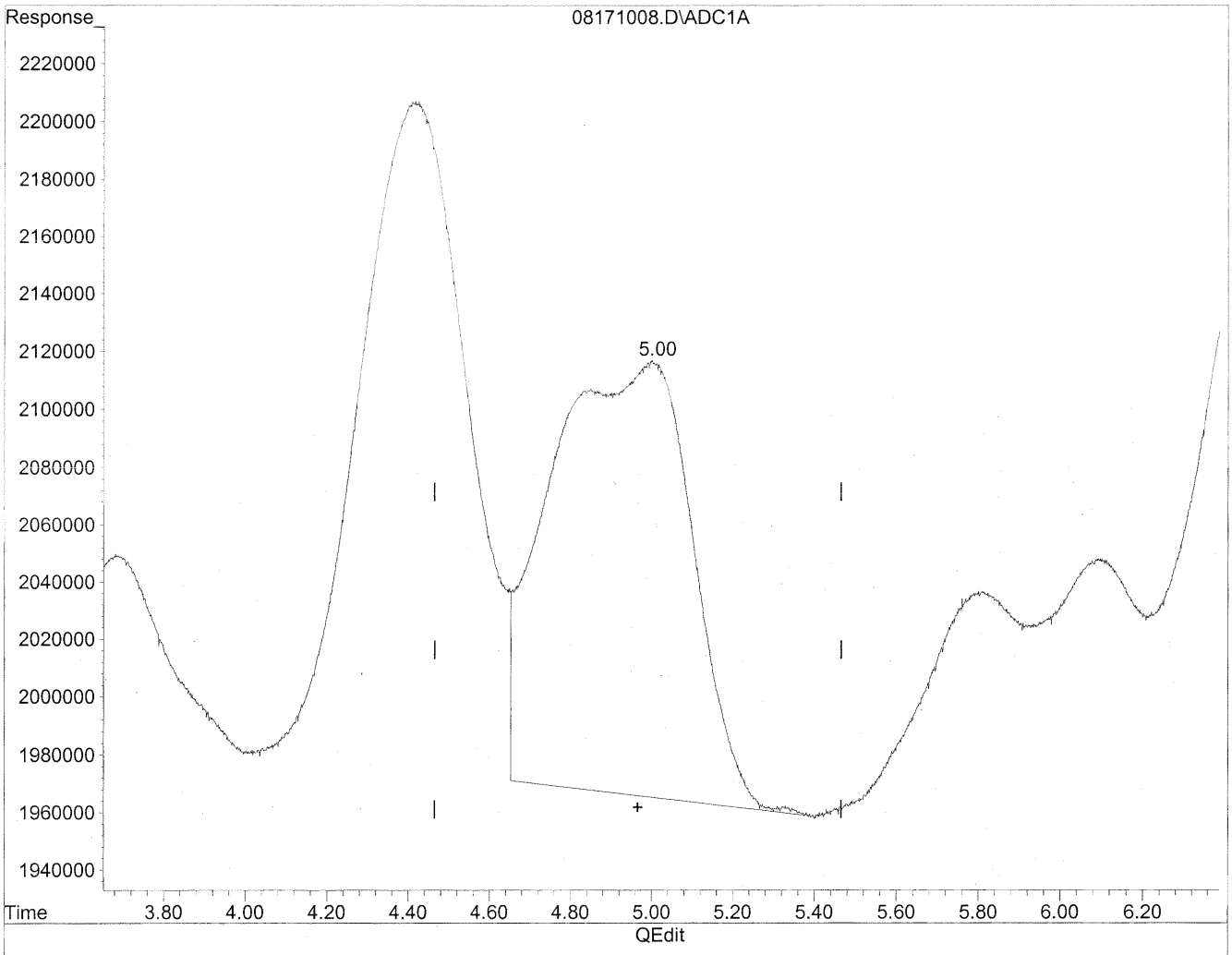
61

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.00min 414.074ng/ml

response 36577666

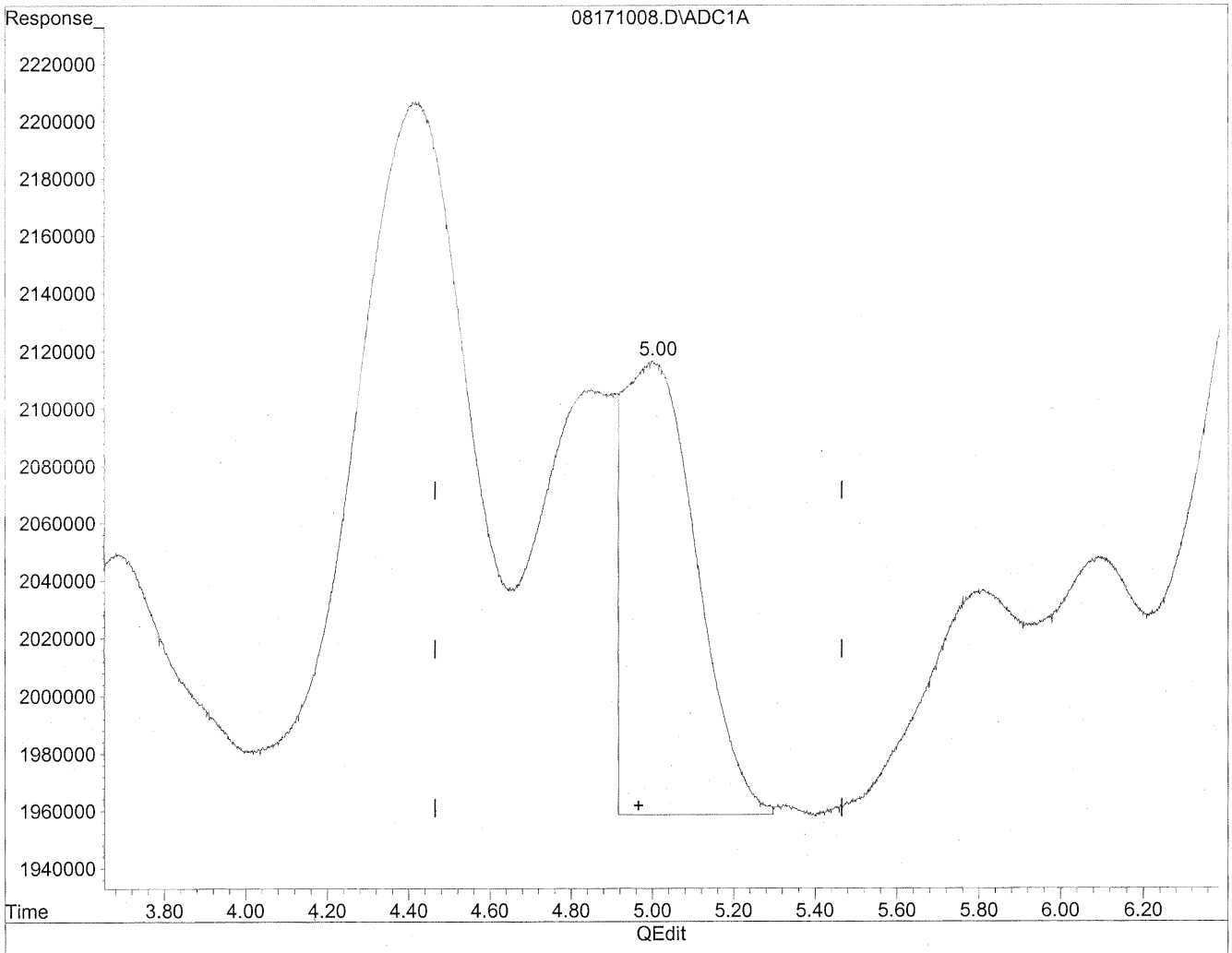
(+) = Expected Retention Time

08171008.D TO110709.M Sat Aug 22 14:13:16 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.00min 223.368ng/ml m

response 19731484

HC 8/22/09

SP

KP 8/23/09

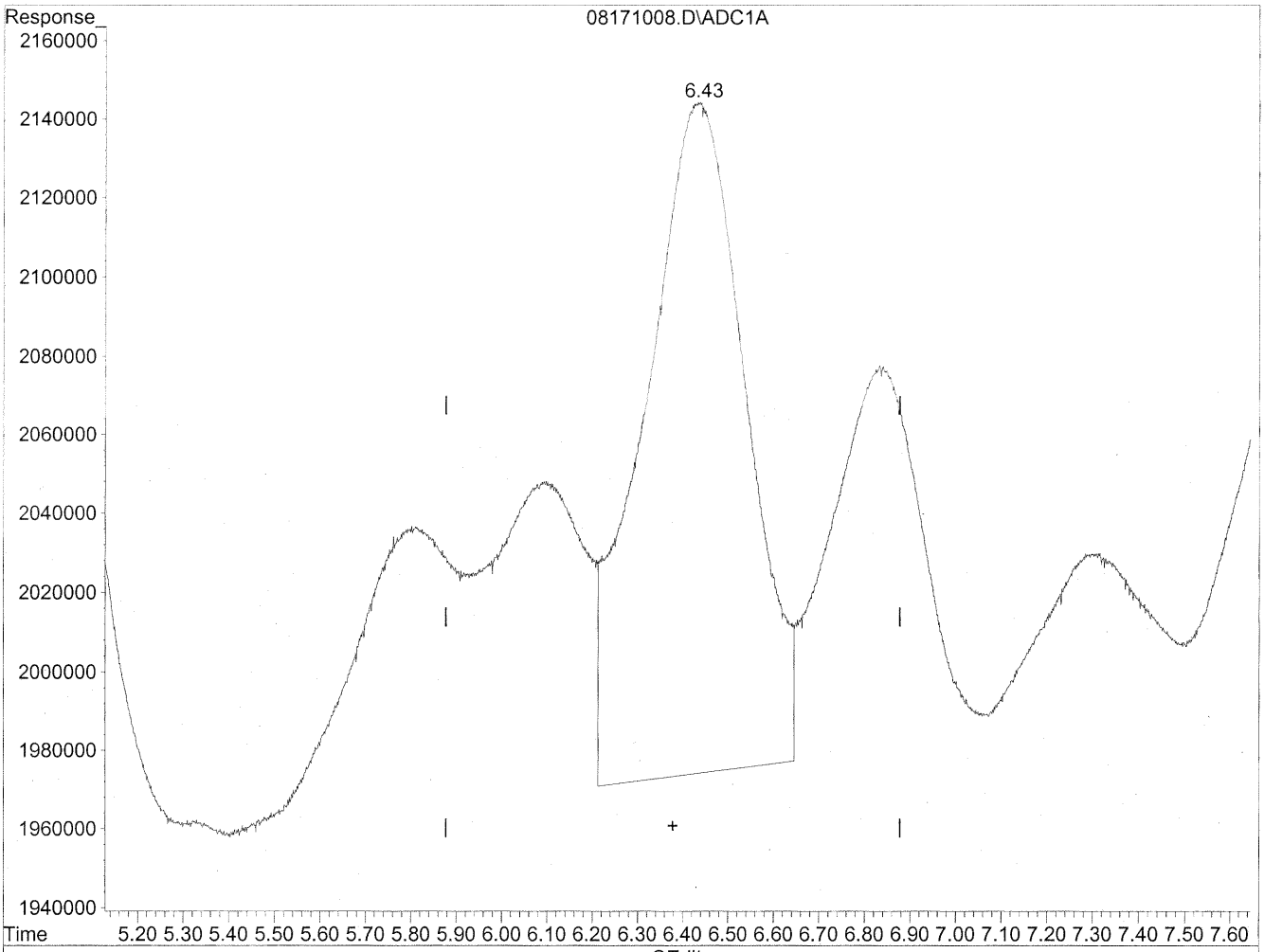
(+) = Expected Retention Time

08171008.D TO110709.M Sat Aug 22 14:13:54 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(6) Benzaldehyde

6.43min 403.049ng/ml

response 26548583

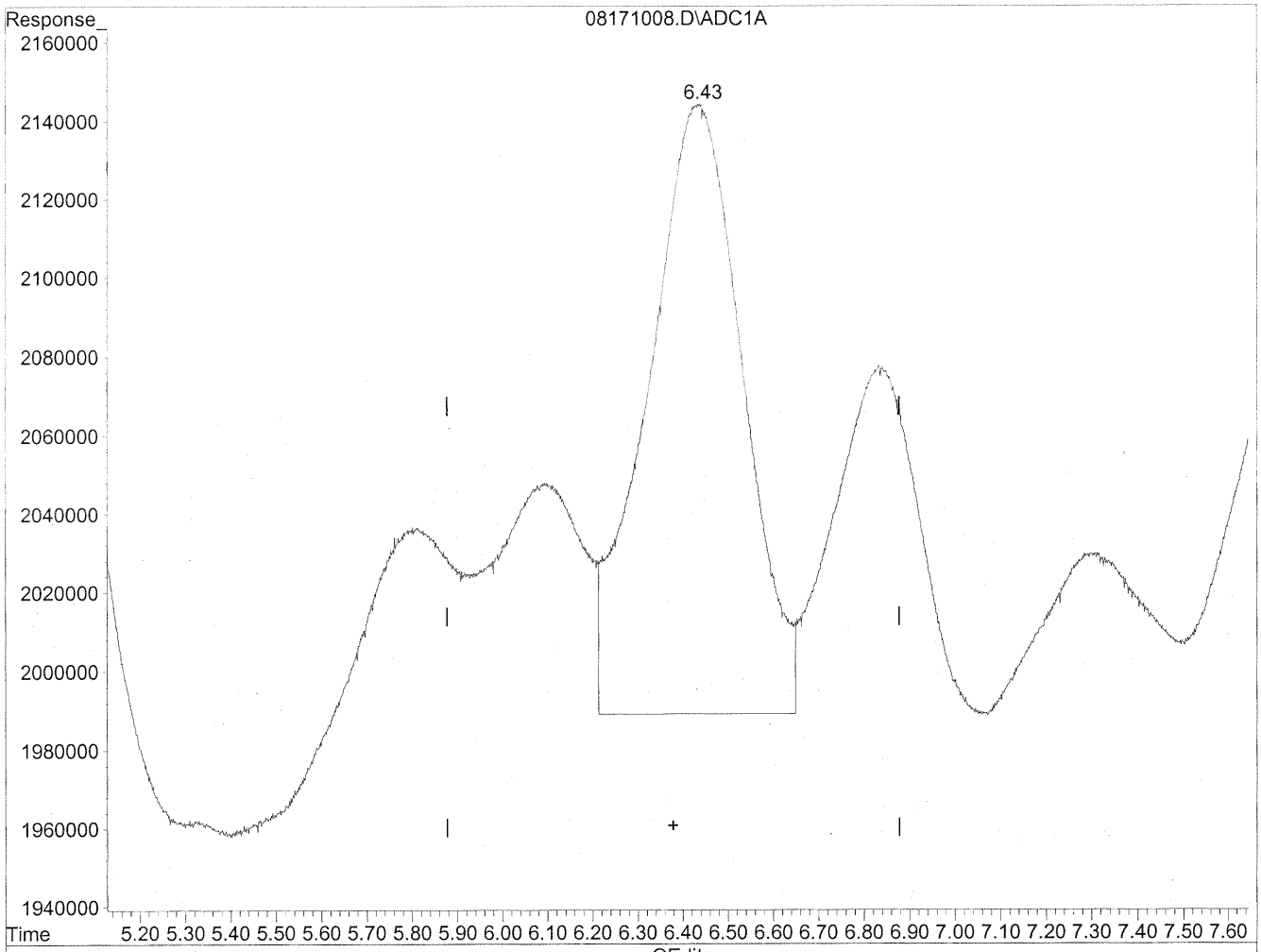
(+) = Expected Retention Time

08171008.D TO110709.M Sat Aug 22 14:13:58 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.43min 344.570ng/ml m

response 22696600

HC 8/22/09
MC
KR 8/23/09

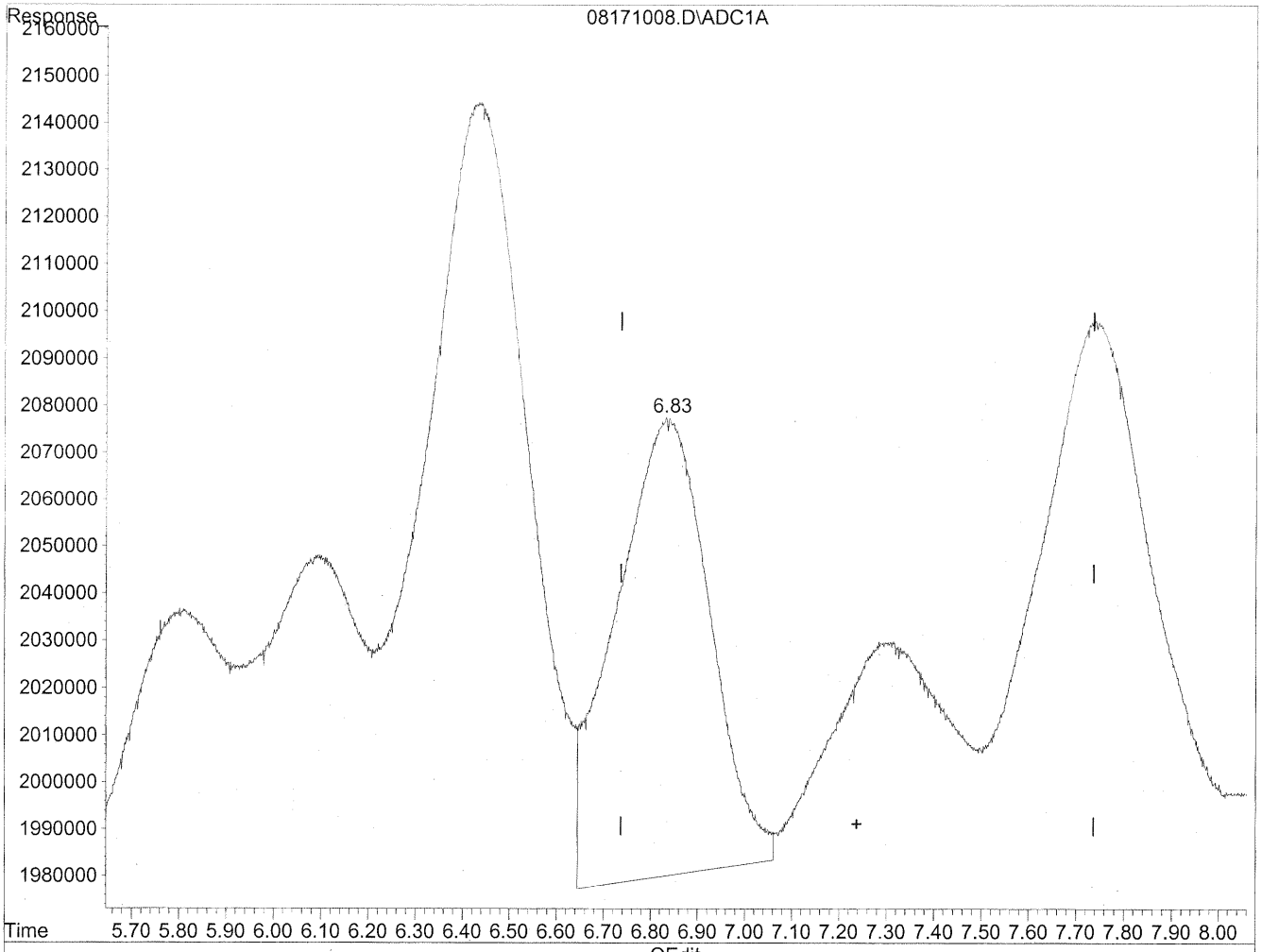
(+) = Expected Retention Time

08171008.D TO110709.M Sat Aug 22 14:14:14 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

6.84min 173.159ng/ml

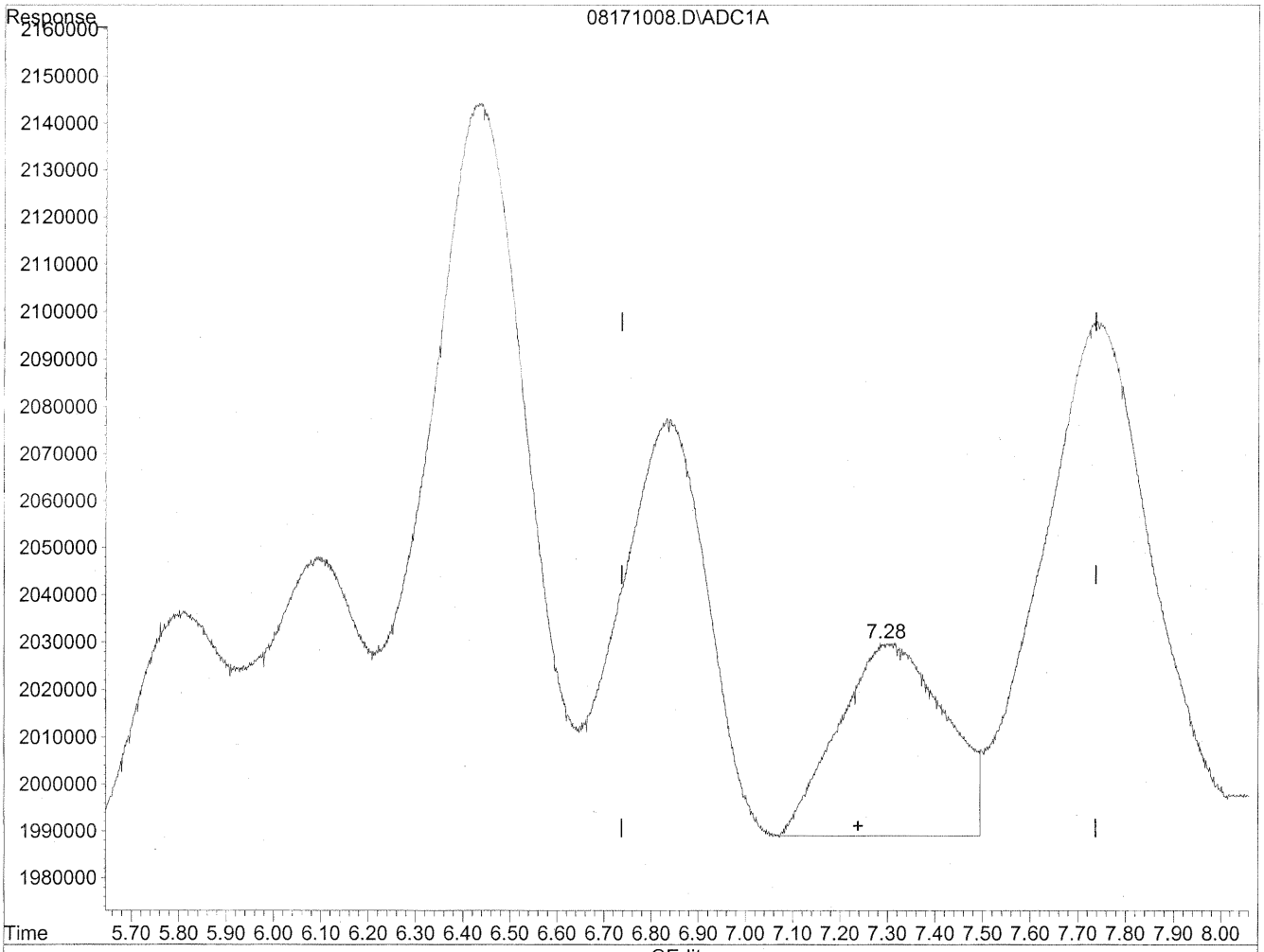
response 13549865

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.28min 81.304ng/ml m

response 6362152

HC 8/22/09 MP

KR 8/23/09

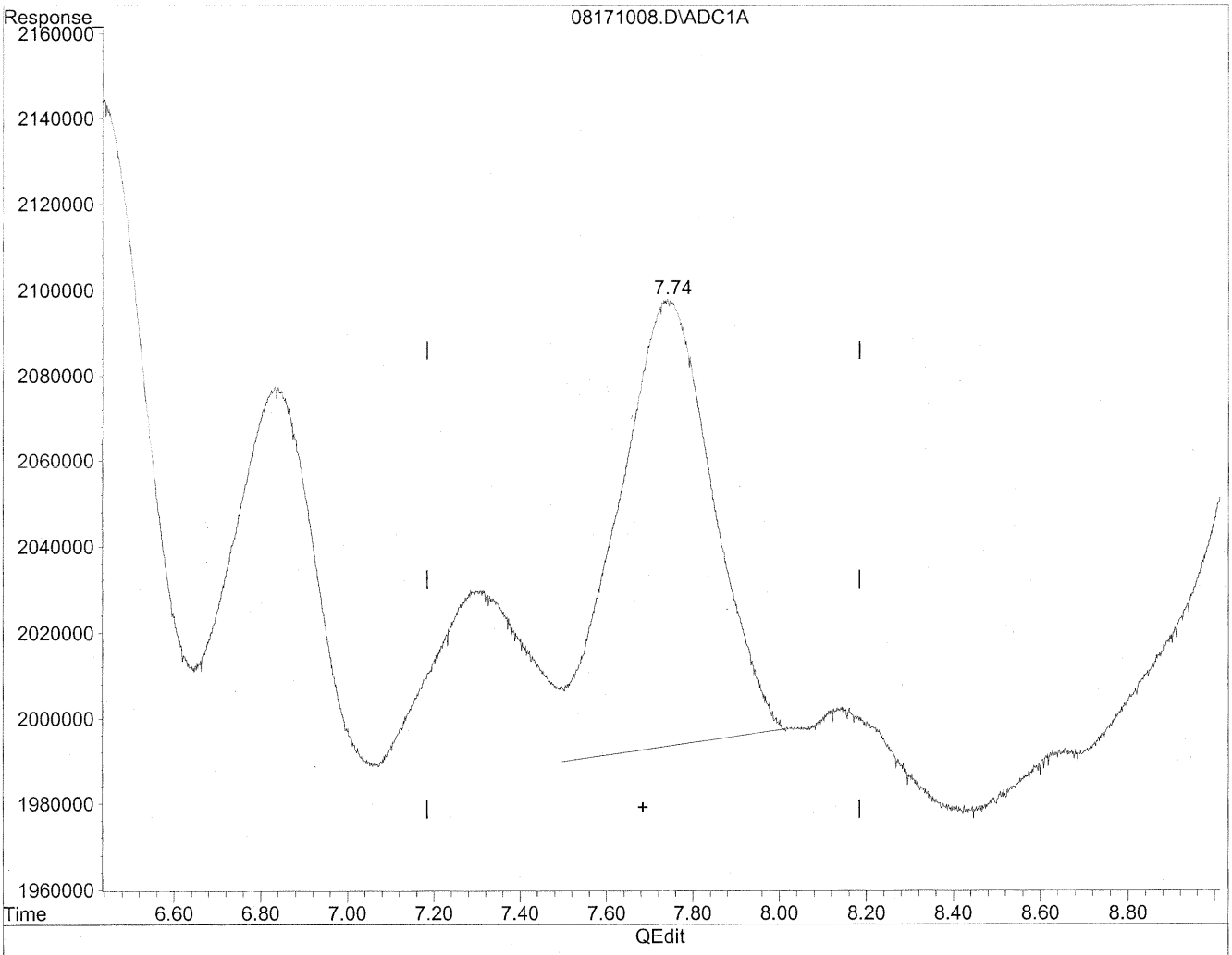
(+) = Expected Retention Time

08171008.D TO110709.M Sat Aug 22 14:14:40 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.74min 218.006ng/ml

response 16024555

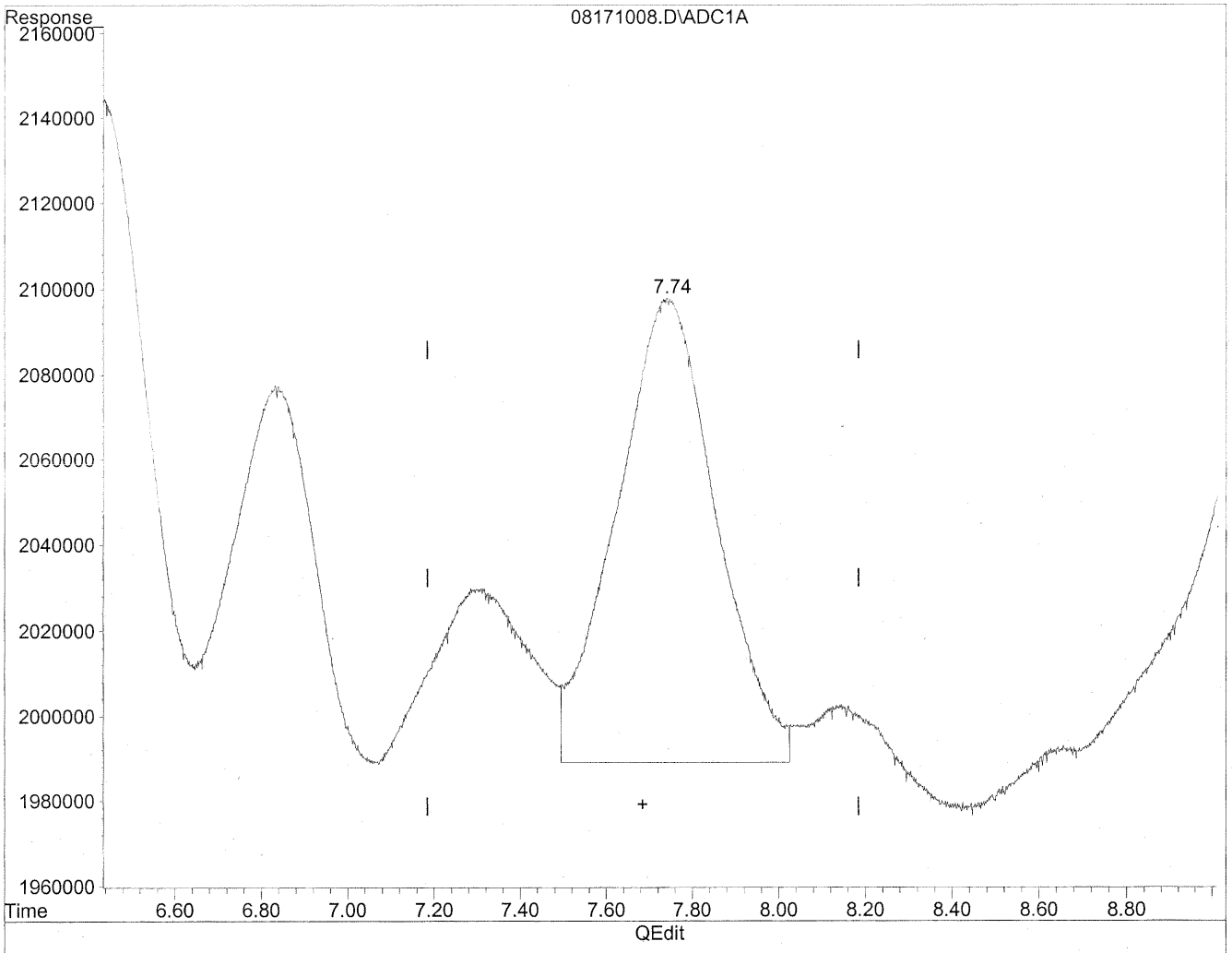
(+) = Expected Retention Time

08171008.D TO110709.M Sat Aug 22 14:14:44 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.74min 238.403ng/ml m

response 17523823

HC 8/22/09 MC

KP 8/23/09

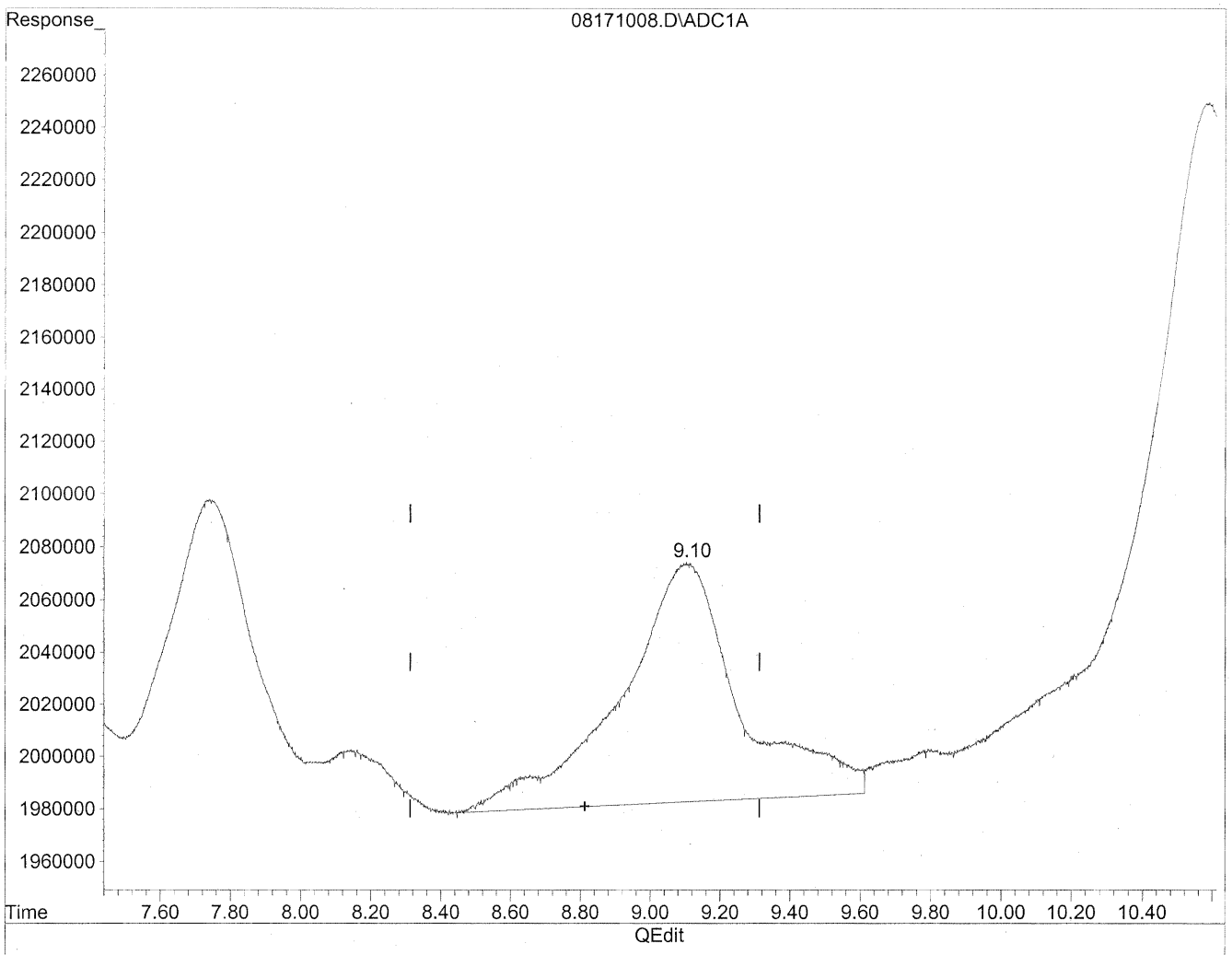
(+) = Expected Retention Time

08171008.D TO110709.M Sat Aug 22 14:15:06 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

9.10min 404.136ng/ml

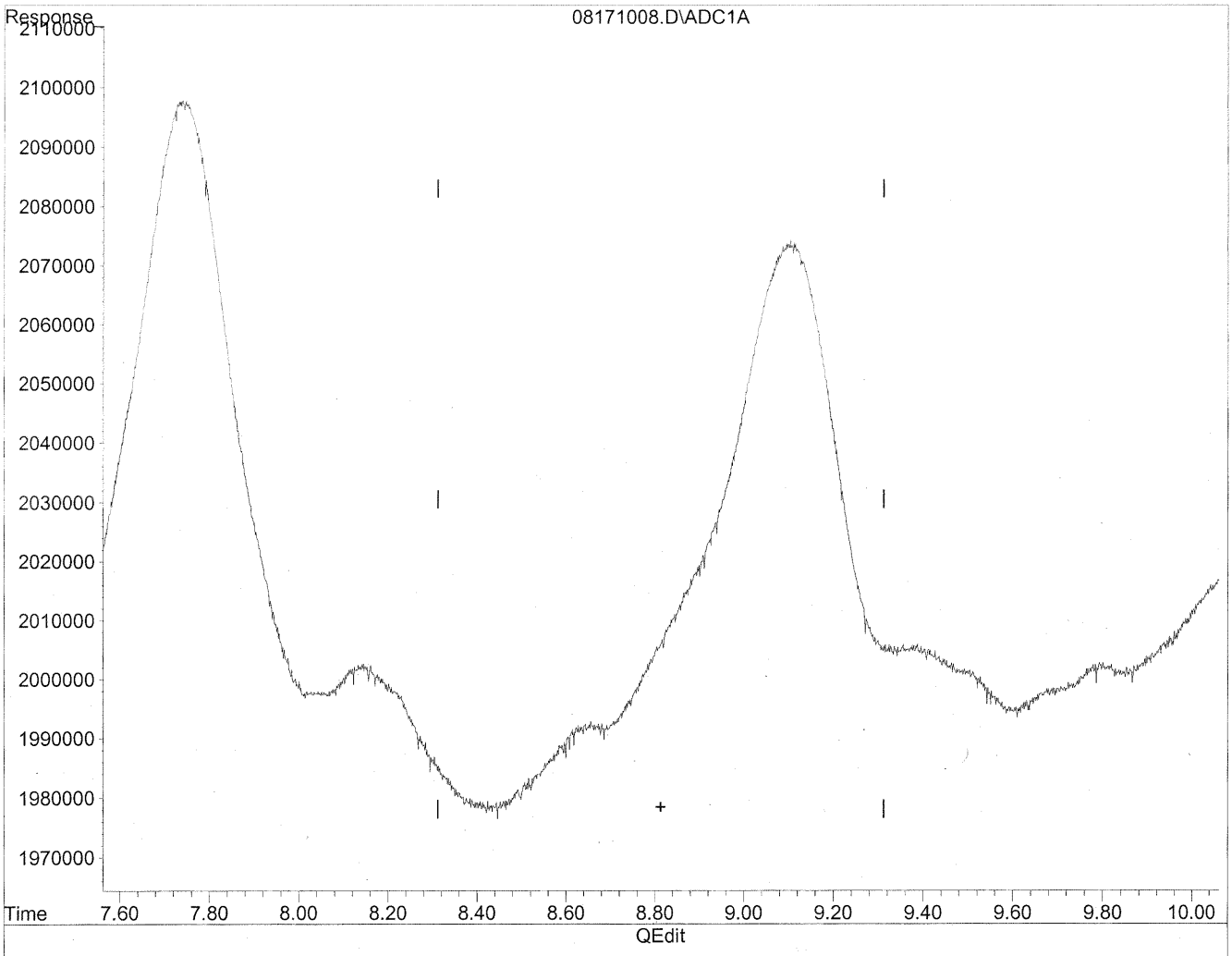
response 21821502

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09 MD

KPC 8/23/09

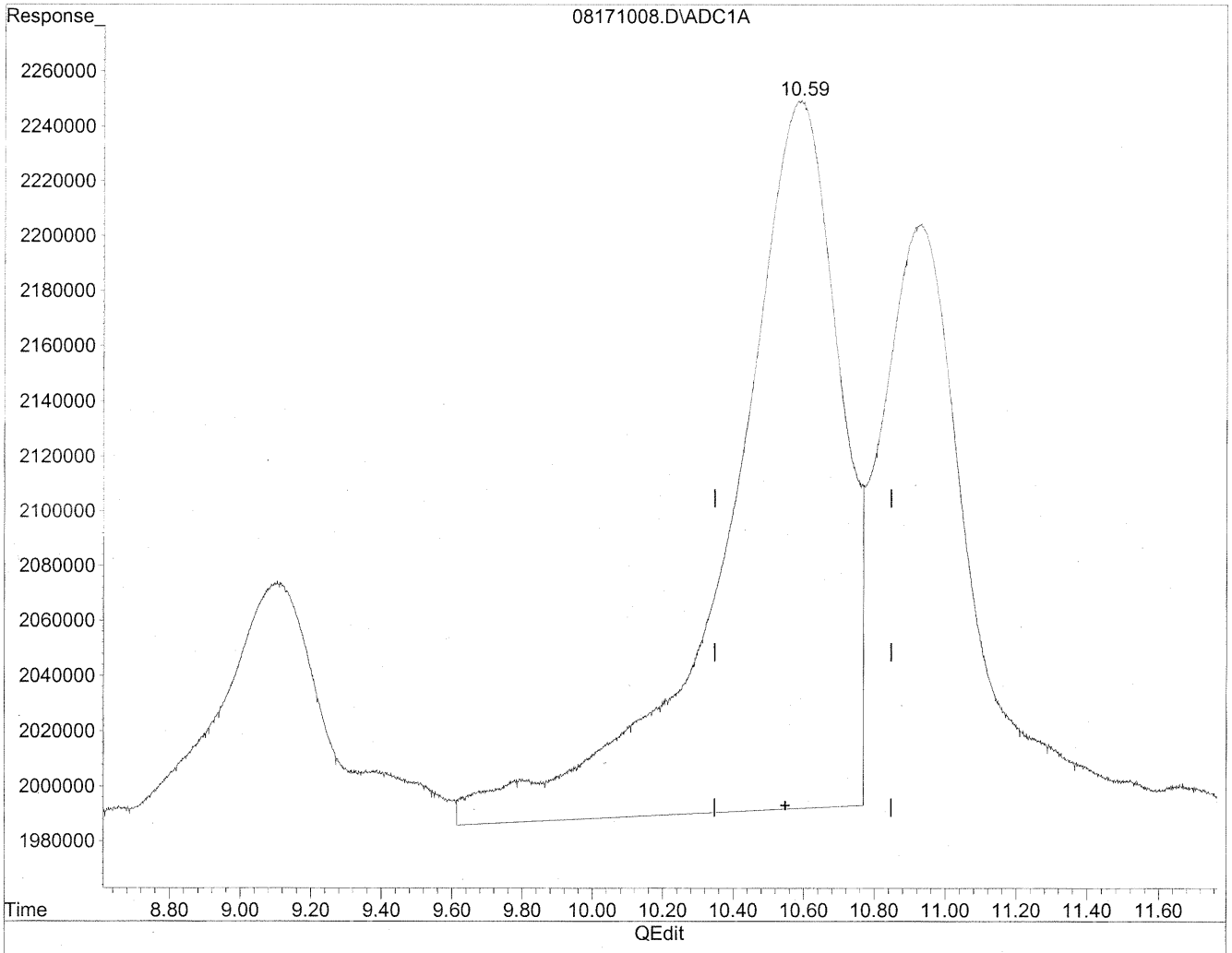
71

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.59min 840.230ng/ml

response 56584243

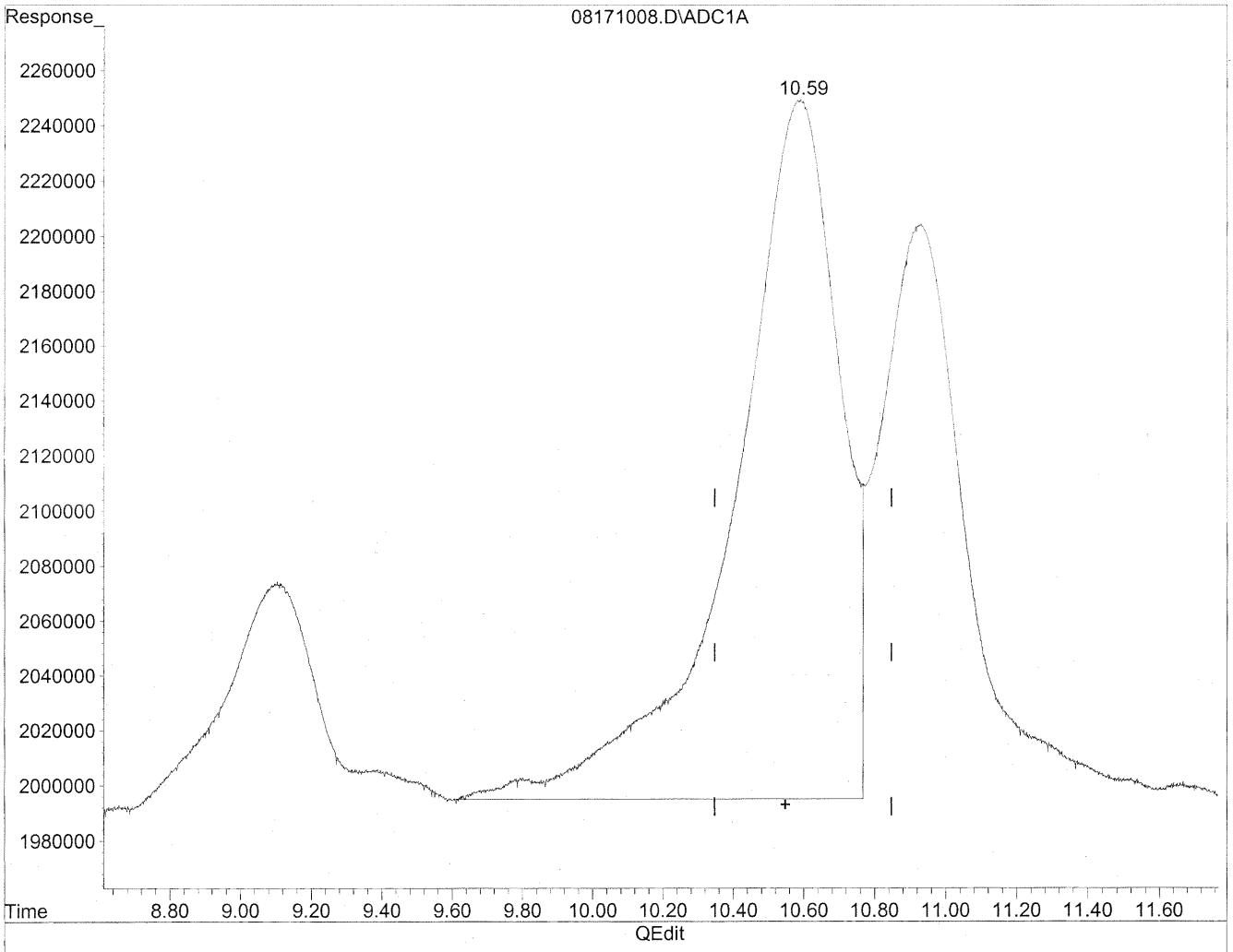
(+) = Expected Retention Time

08171008.D TO110709.M Sat Aug 22 14:15:20 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.59min 781.477ng/ml m

response 52627599

tlc 8/22/09 IC

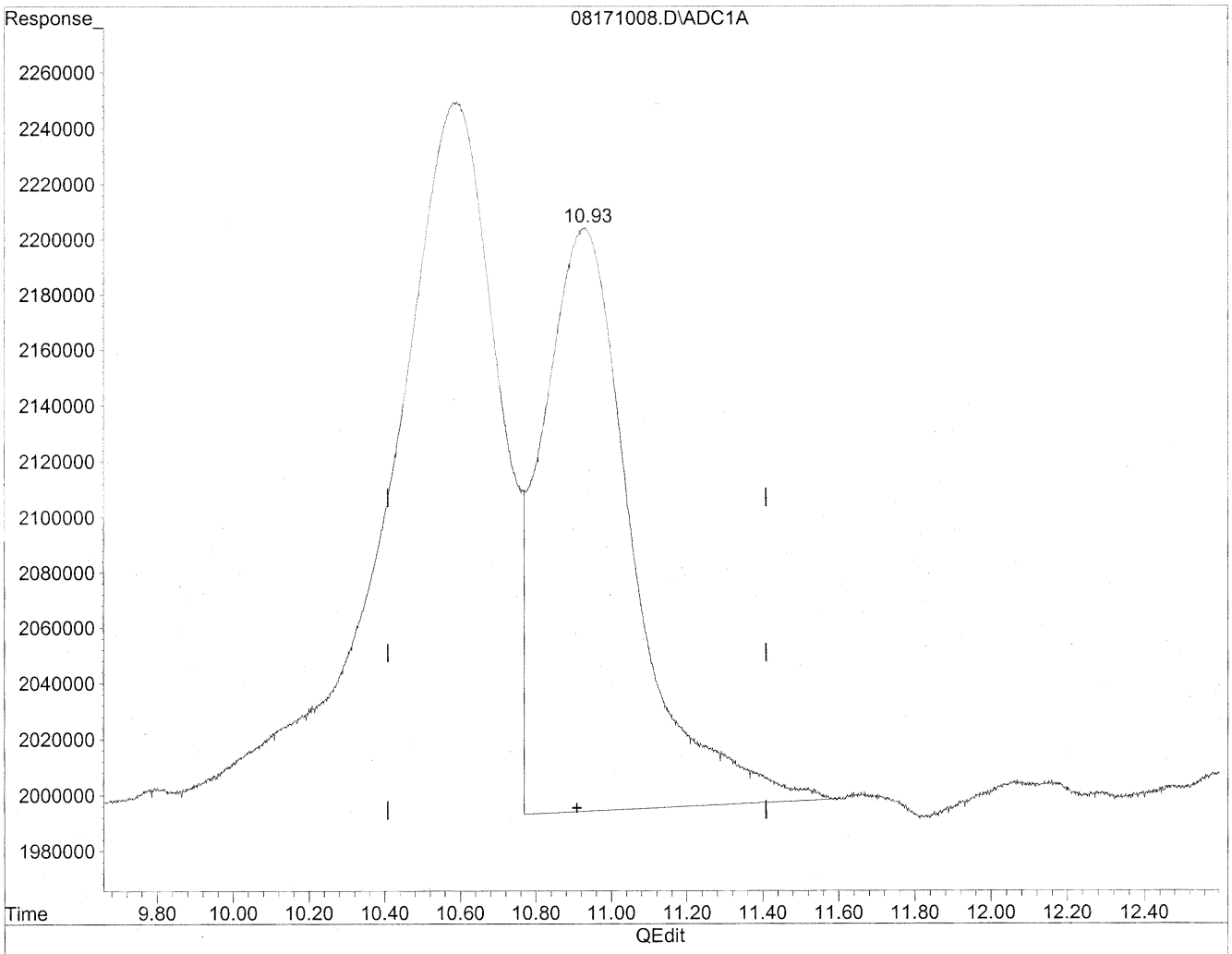
KP 8/23/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.93min 713.460ng/ml

response 34969104

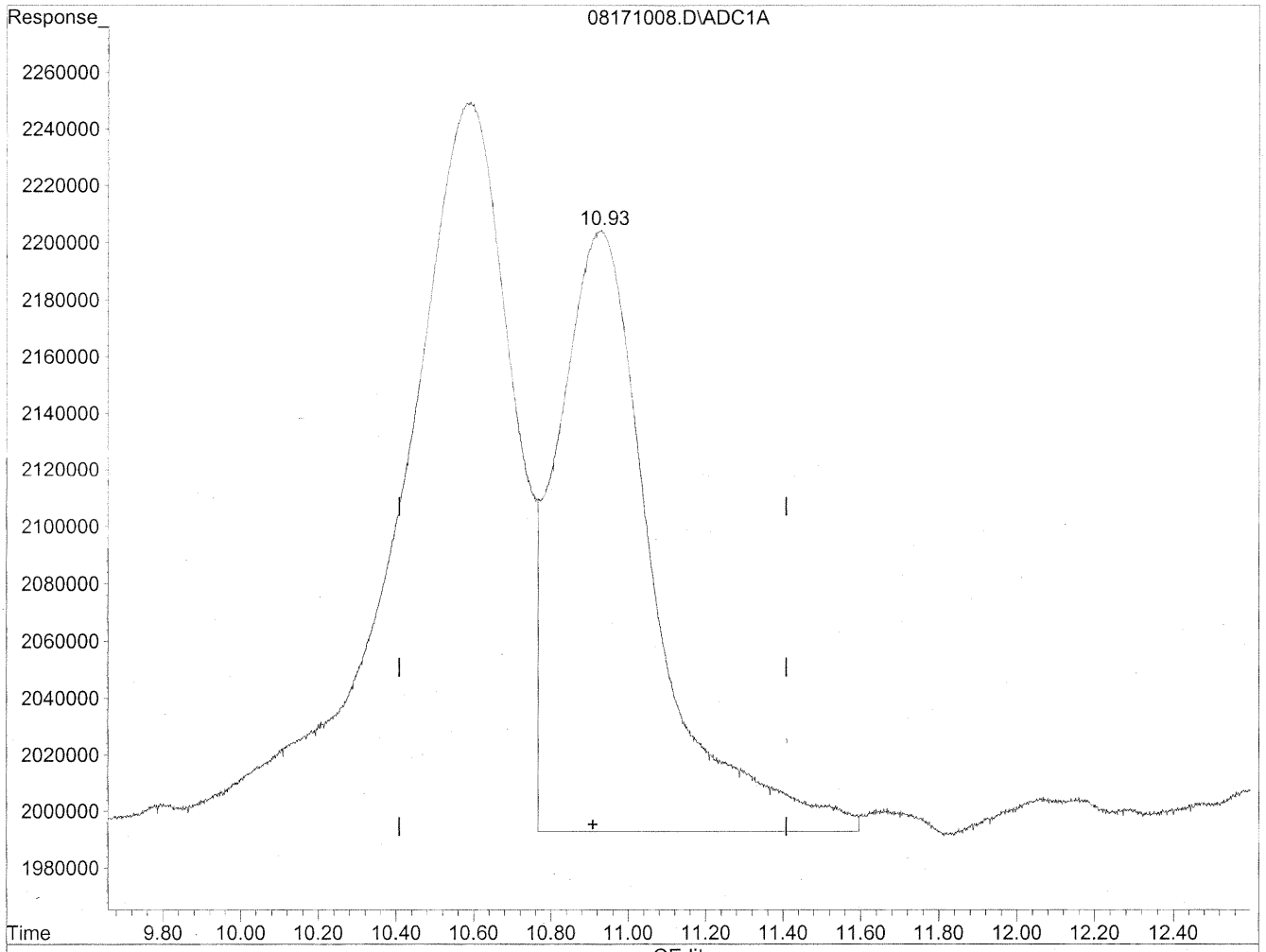
(+) = Expected Retention Time

08171008.D TO110709.M Sat Aug 22 14:15:34 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171008.D Vial: 21
Acq On : 18 Aug 2009 5:39 pm Operator: HC
Sample : P0902772-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:11 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.93min 744.633ng/ml m

response 36497004

HC 8/22/09 LC

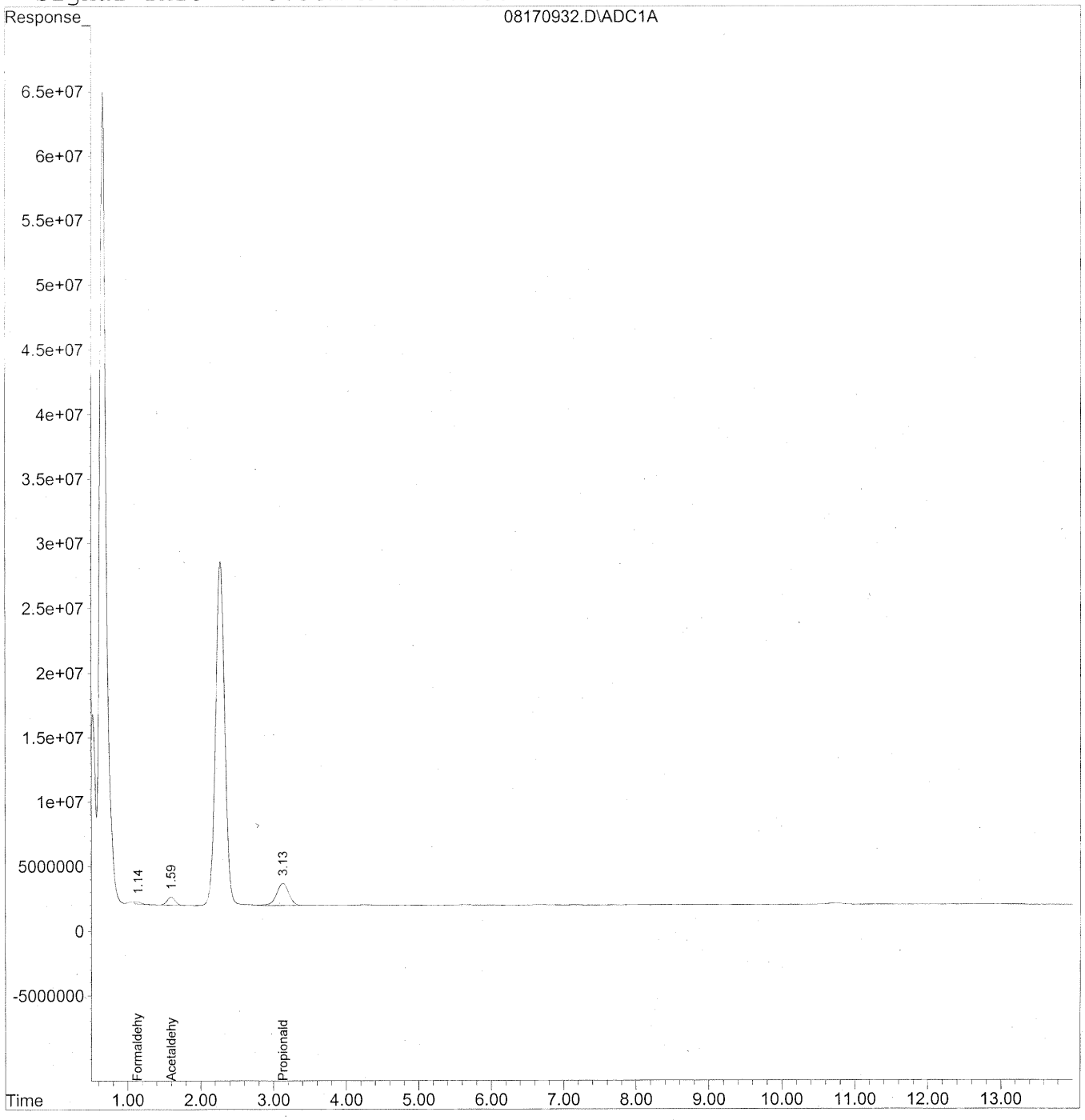
KP 8/23/09

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_08\17\08170932.D Vial: 31
Acq On : 17 Aug 2009 10:36 pm Operator: HC
Sample : P0902772-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:41 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 17 18:29:01 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170932.D Vial: 31
Acq On : 17 Aug 2009 10:36 pm Operator: HC
Sample : P0902772-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:41 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 17 18:29:01 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.13	10066094	54.832	ng/ml
2)	Acetaldehyde	1.59	48171651	343.535	ng/ml
3)	Propionaldehyde	3.13	200867563	~~1882.629~~	~~ng/ml~~
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/mld
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/mld

HL 8/28/09

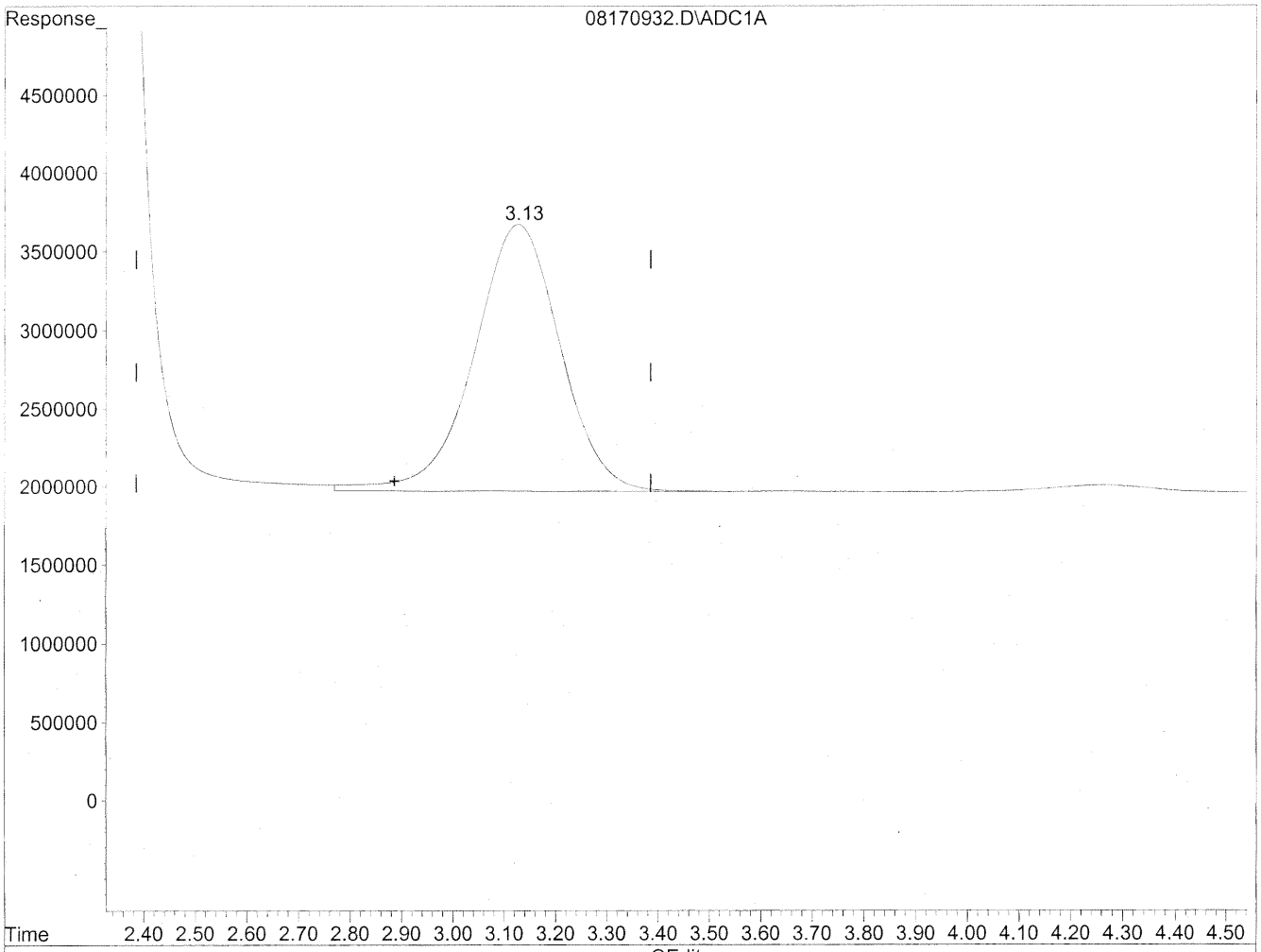
77

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170932.D Vial: 31
Acq On : 17 Aug 2009 10:36 pm Operator: HC
Sample : P0902772-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:41 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.13min 1882.629ng/ml

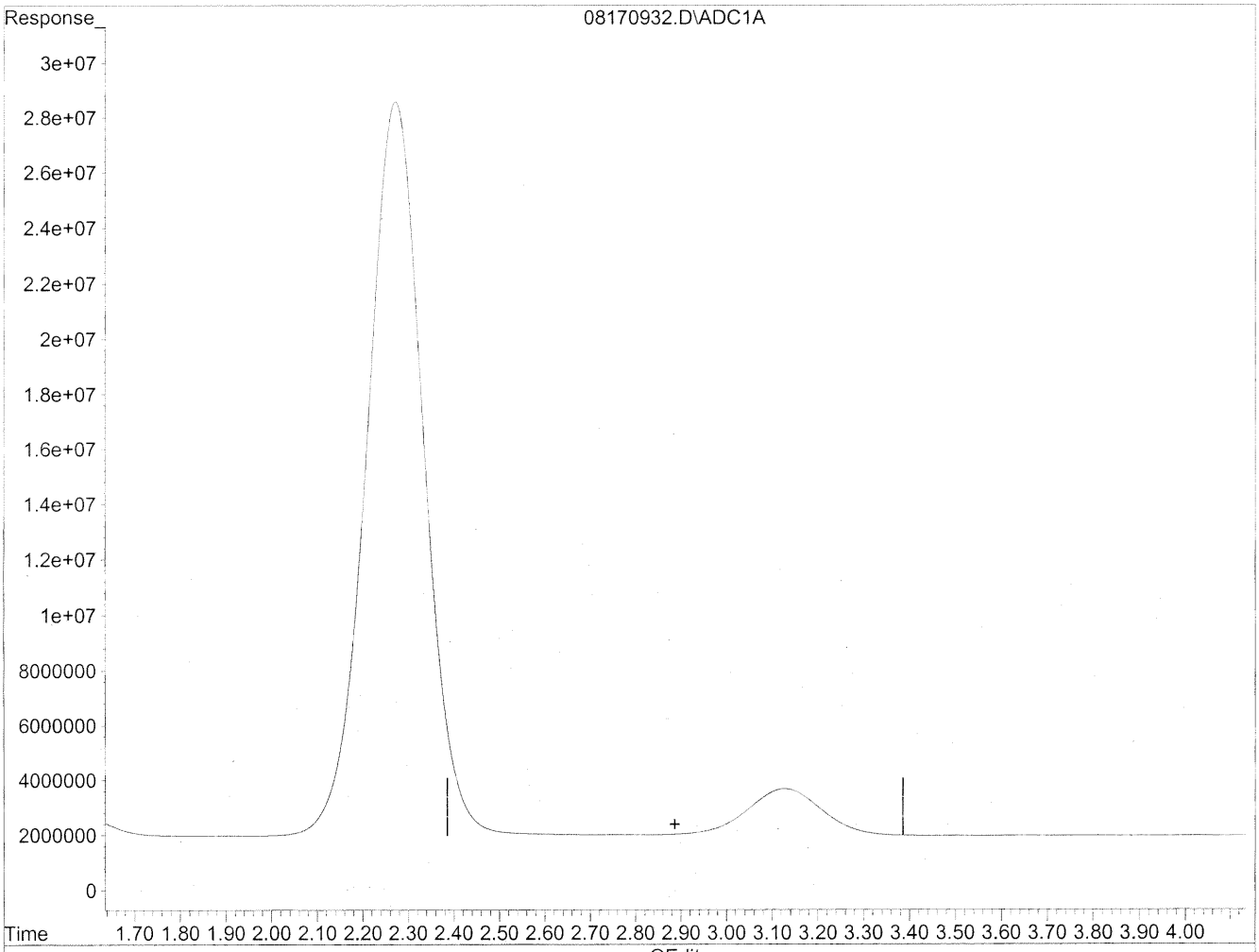
response 200867563

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170932.D Vial: 31
Acq On : 17 Aug 2009 10:36 pm Operator: HC
Sample : P0902772-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:41 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

tic 8/25/09 MP

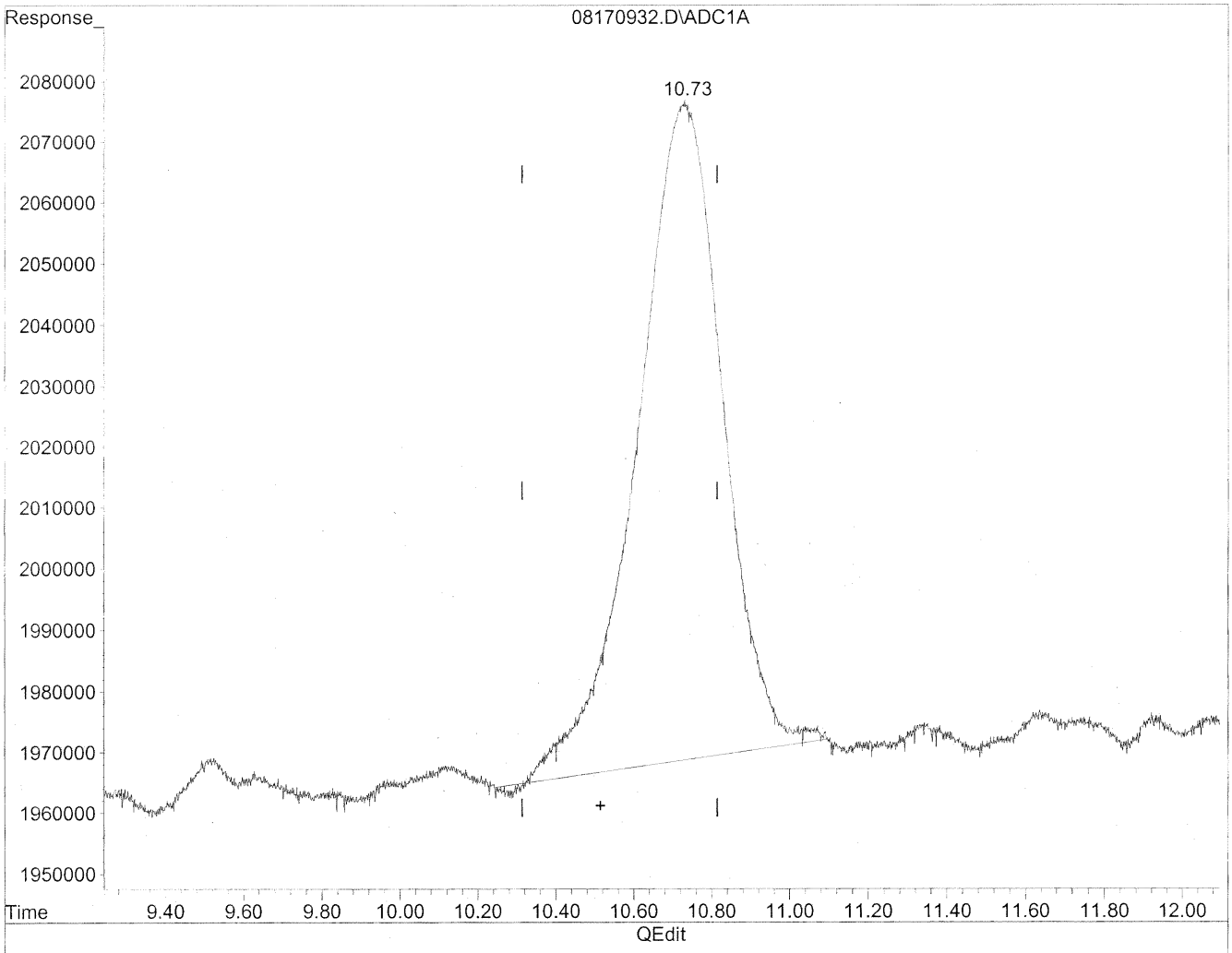
KB 8/25/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170932.D Vial: 31
Acq On : 17 Aug 2009 10:36 pm Operator: HC
Sample : P0902772-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 18 11:25 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.73min 243.755ng/ml

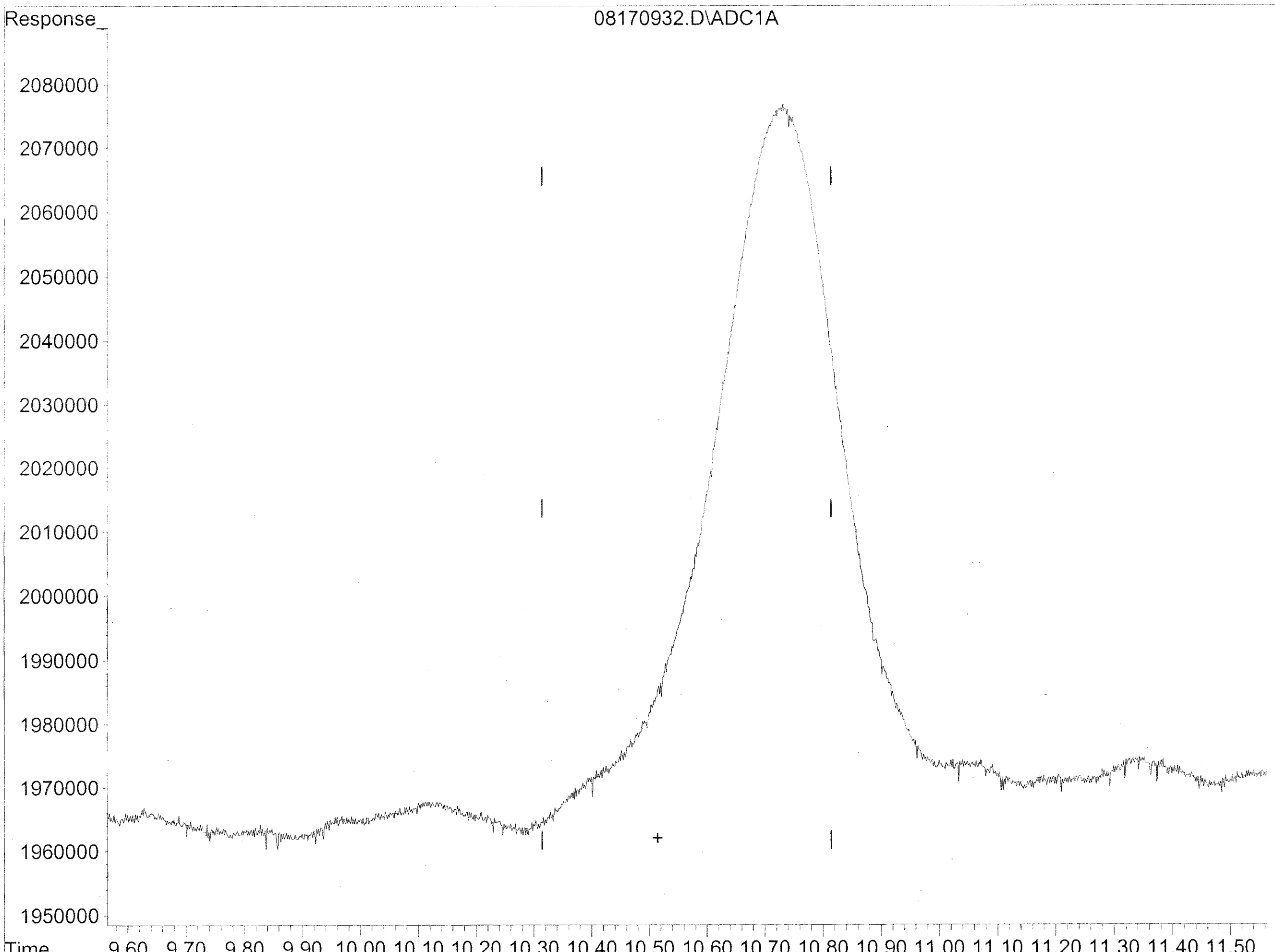
response 16415402

(+) = Expected Retention Time

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2009_08\17\08170932.D        Vial: 31
Acq On    : 17 Aug 2009   10:36 pm                      Operator: HC
Sample    : P0902772-003 back 1.0ml                     Inst    : LC 01
Misc      :                                             Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Aug 18 11:25 19109  Quant Results File: TO110709.RES

Method         : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title          : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update    : Mon Aug 24 08:44:34 2009
Response via   : Multiple Level Calibration
```



QEdit

(11) Hexaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/25/09 up

KR 8/25/09

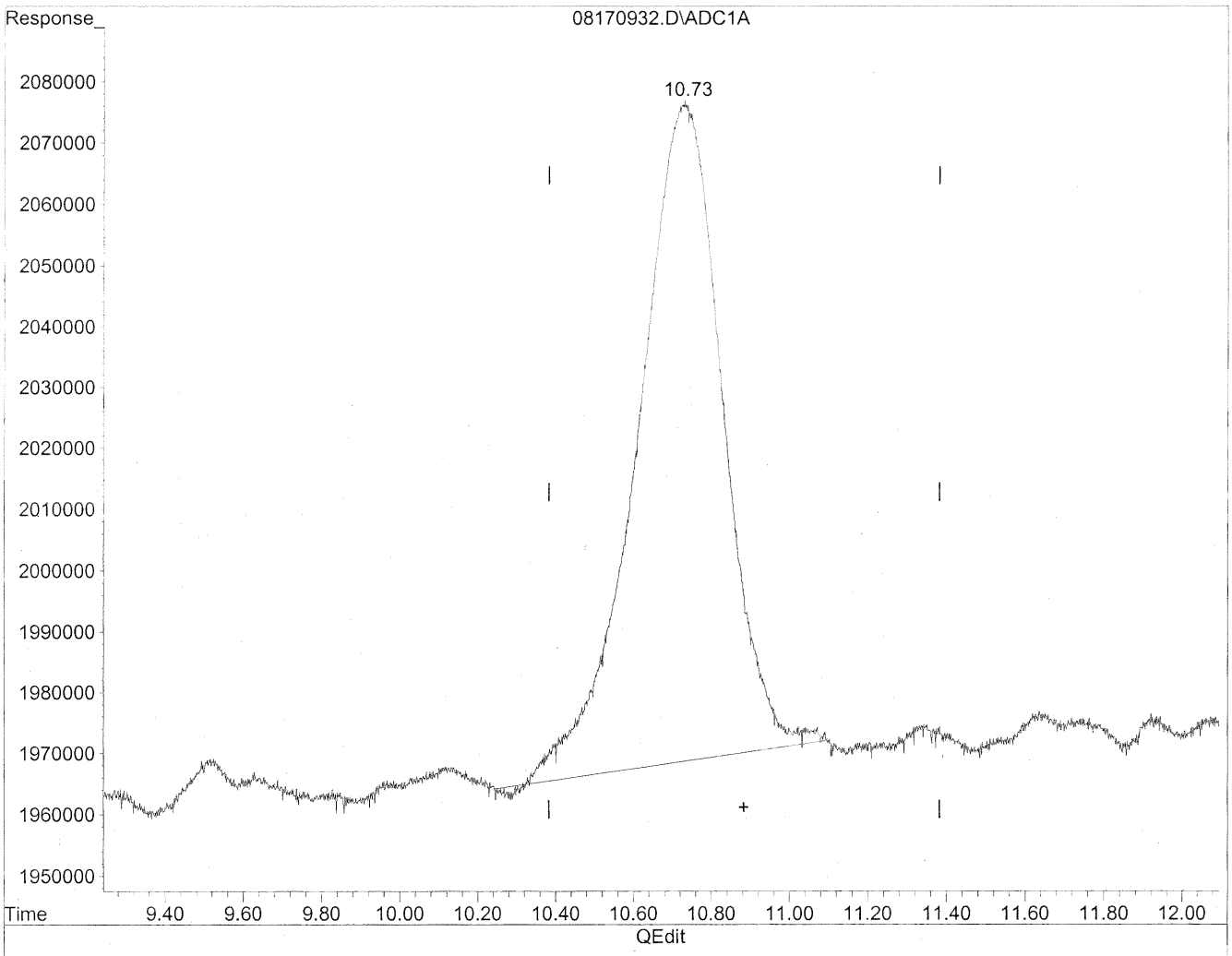
(+) = Expected Retention Time

08170932.D TO110709.M Tue Aug 25 07:40:38 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170932.D Vial: 31
Acq On : 17 Aug 2009 10:36 pm Operator: HC
Sample : P0902772-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 18 11:25 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.73min 334.917ng/ml

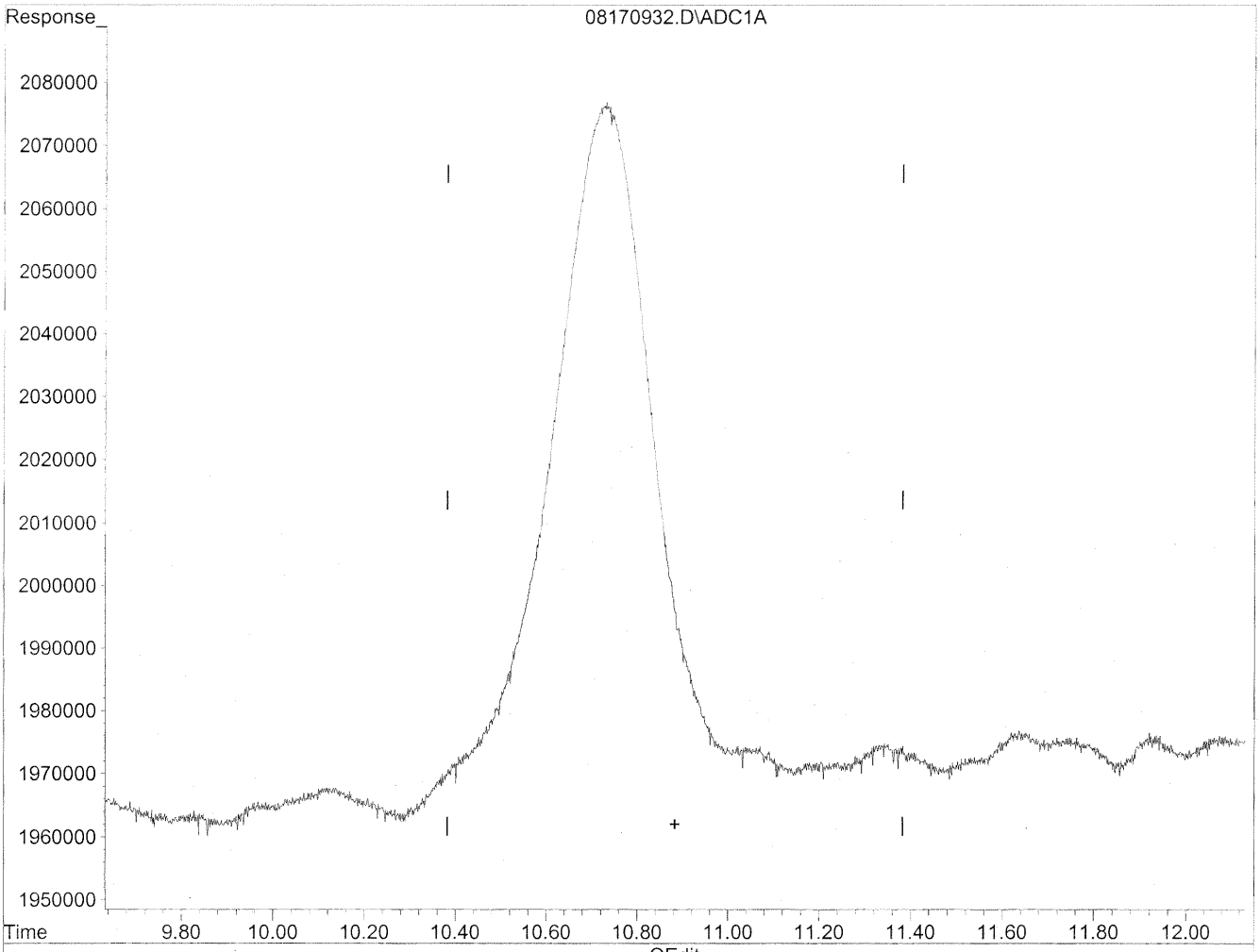
response 16415402

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170932.D Vial: 31
Acq On : 17 Aug 2009 10:36 pm Operator: HC
Sample : P0902772-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 18 11:25 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

0.00min 0.000ng/ml d

response 0

ll 8/25/09 mf

KLR 8/25/09

(+) = Expected Retention Time

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **100306**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P0902772-004

Test Code:	EPA Method TO-11A	Date Collected:	8/10/09
Instrument ID:	Waters LC Module I Plus/UV_Vis 360/LC1	Date Received:	8/12/09
Analyst:	Hani Cherazaie	Date Analyzed:	8/17 - 8/18/09
Sampling Media:	Silica Gel DNPH Tube	Desorption Volume:	1.0 ml
Test Notes:	**BC**	Volume Sampled:	109.6 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**4,200**	**38**	0.91	**31**	0.74	
75-07-0	**Acetaldehyde**	**2,700**	**24**	0.91	**13**	0.51	**BT**
123-38-6	**Propionaldehyde**	**220**	**2.0**	0.91	**0.85**	0.38	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.91	ND	0.32	
123-72-8	**Butyraldehyde**	**240**	**2.2**	0.91	**0.74**	0.31	
100-52-7	**Benzaldehyde**	**350**	**3.2**	0.91	**0.74**	0.21	
590-86-3	Isovaleraldehyde	< 100	ND	0.91	ND	0.26	
110-62-3	**Valeraldehyde**	**330**	**3.0**	0.91	**0.86**	0.26	
529-20-4	o-Tolualdehyde	< 100	ND	0.91	ND	0.19	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	1.8	ND	0.37	
66-25-1	**n-Hexaldehyde**	**670**	**6.1**	0.91	**1.5**	0.22	
5779-94-2	**2,5-Dimethylbenzaldehyde**	**580**	**5.3**	0.91	**0.96**	0.17	

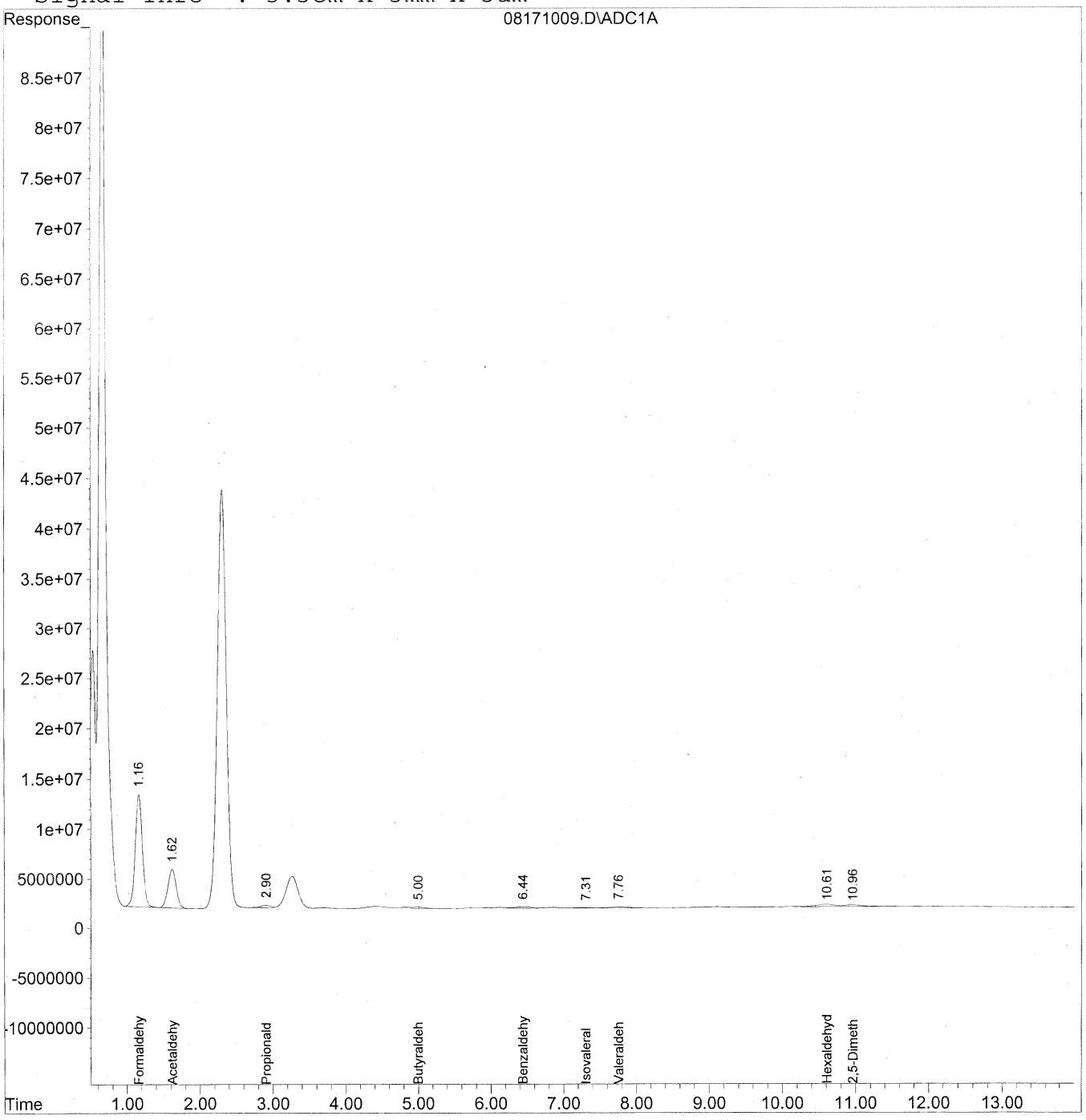
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
BT = Results indicated possible breakthrough; back section > 10% front section.

Verified By: _____ Date: 8/25/09

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:21 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:21 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

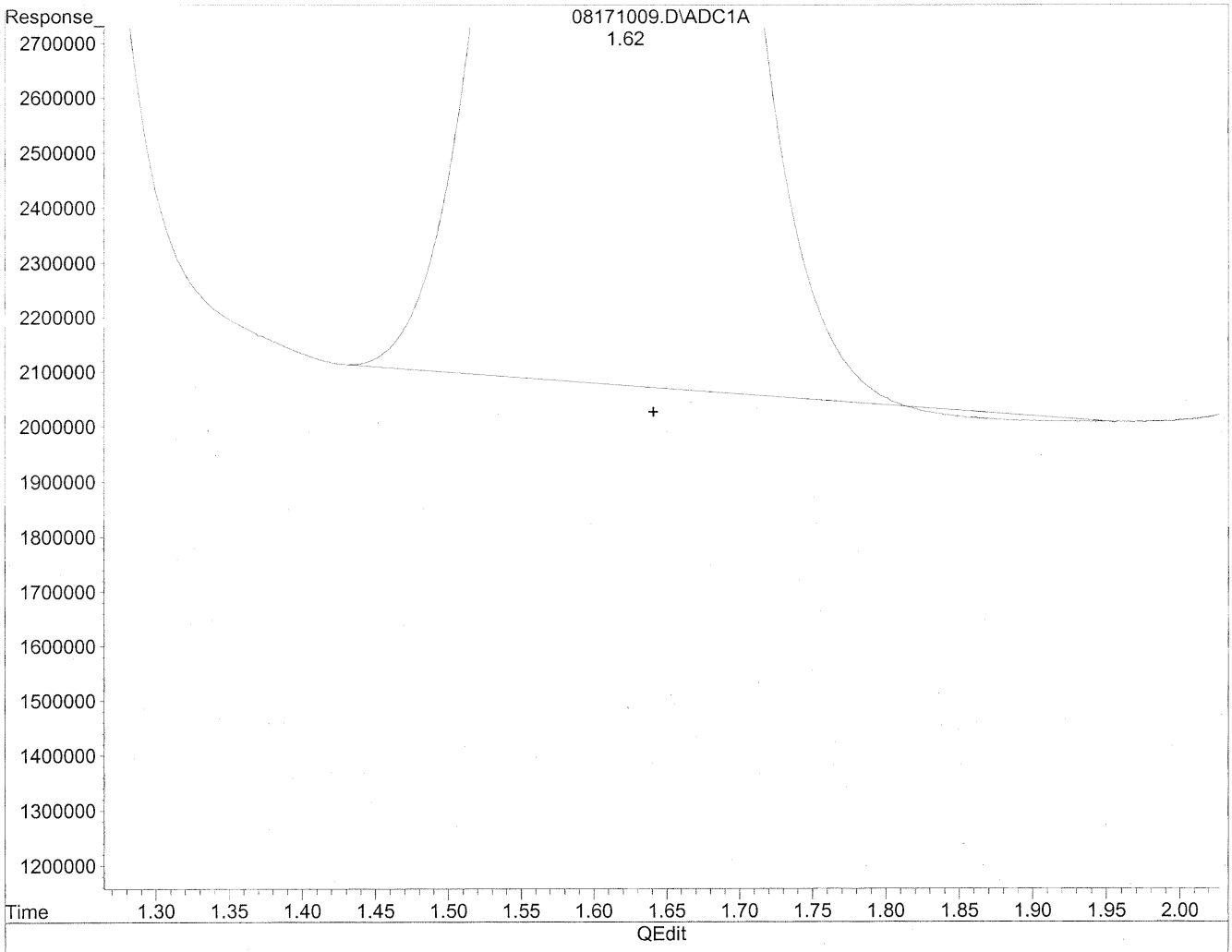
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	763121431	4156.856	ng/ml
2)	Acetaldehyde	1.62	313793211	2237.807	ng/mlm
3)	Propionaldehyde	2.90	23498130	220.236	ng/mlm
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	5.00	21256004	240.626	ng/mlm
6)	Benzaldehyde	6.44	23208609	352.343	ng/mlm
7)	Isovaleraldehyde	7.31	6013814	76.853	ng/mlm
8)	Valeraldehyde	7.76	24404666	332.014	ng/mlm
9)	o-Tolualdehyde	0.00	0	N.D.	ng/mld
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/mld
11)	Hexaldehyde	10.61	45086096	669.492	ng/mlm
12)	2,5-Dimethylbenzaldehyde	10.95	28308687	577.570	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.62min 2223.764ng/ml

response 311824048

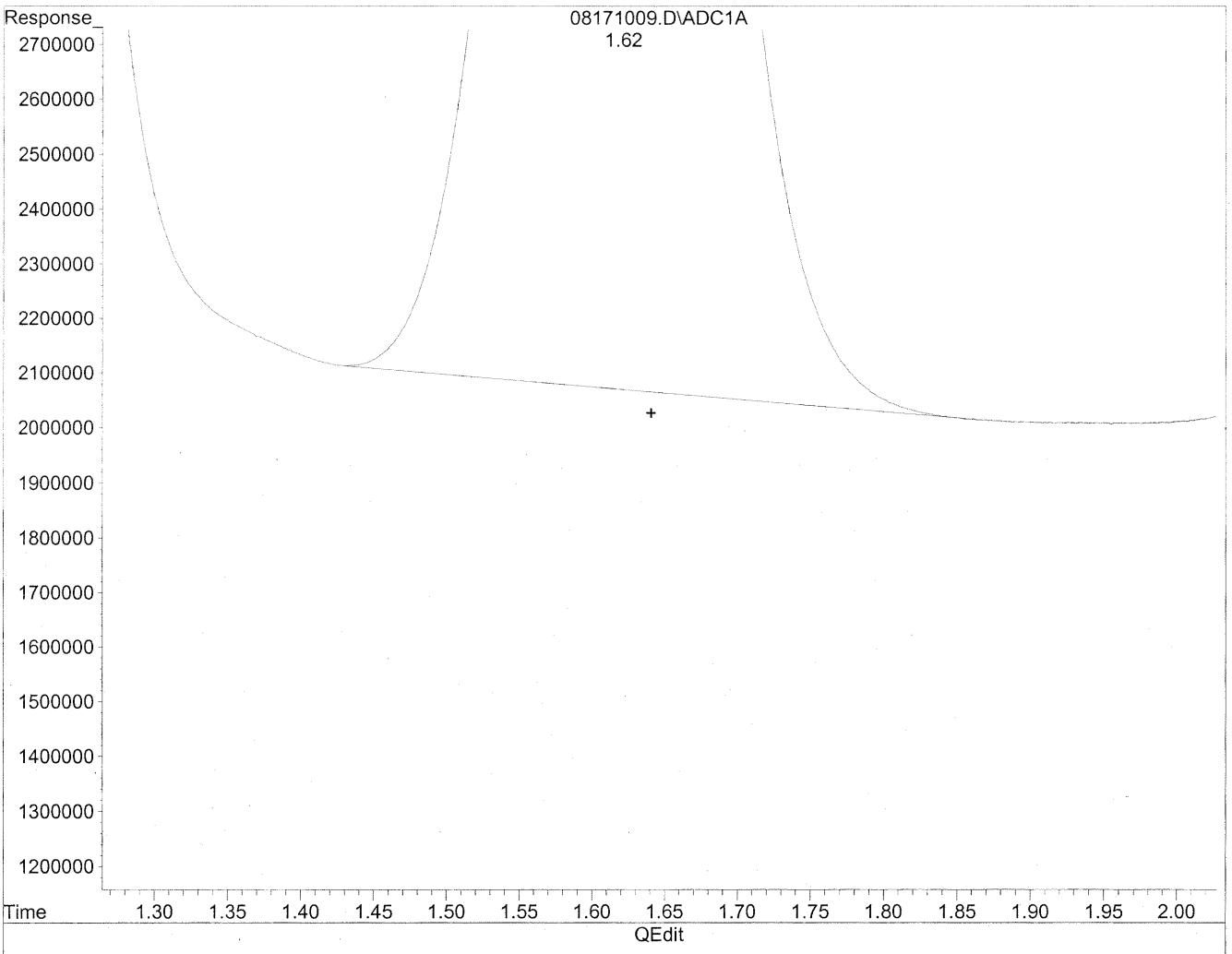
(+) = Expected Retention Time

08171009.D TO110709.M Sat Aug 22 14:18:46 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.62min 2237.807ng/ml m

response 313793211

HC 8/22/09 1C

KP 8/23/09

88

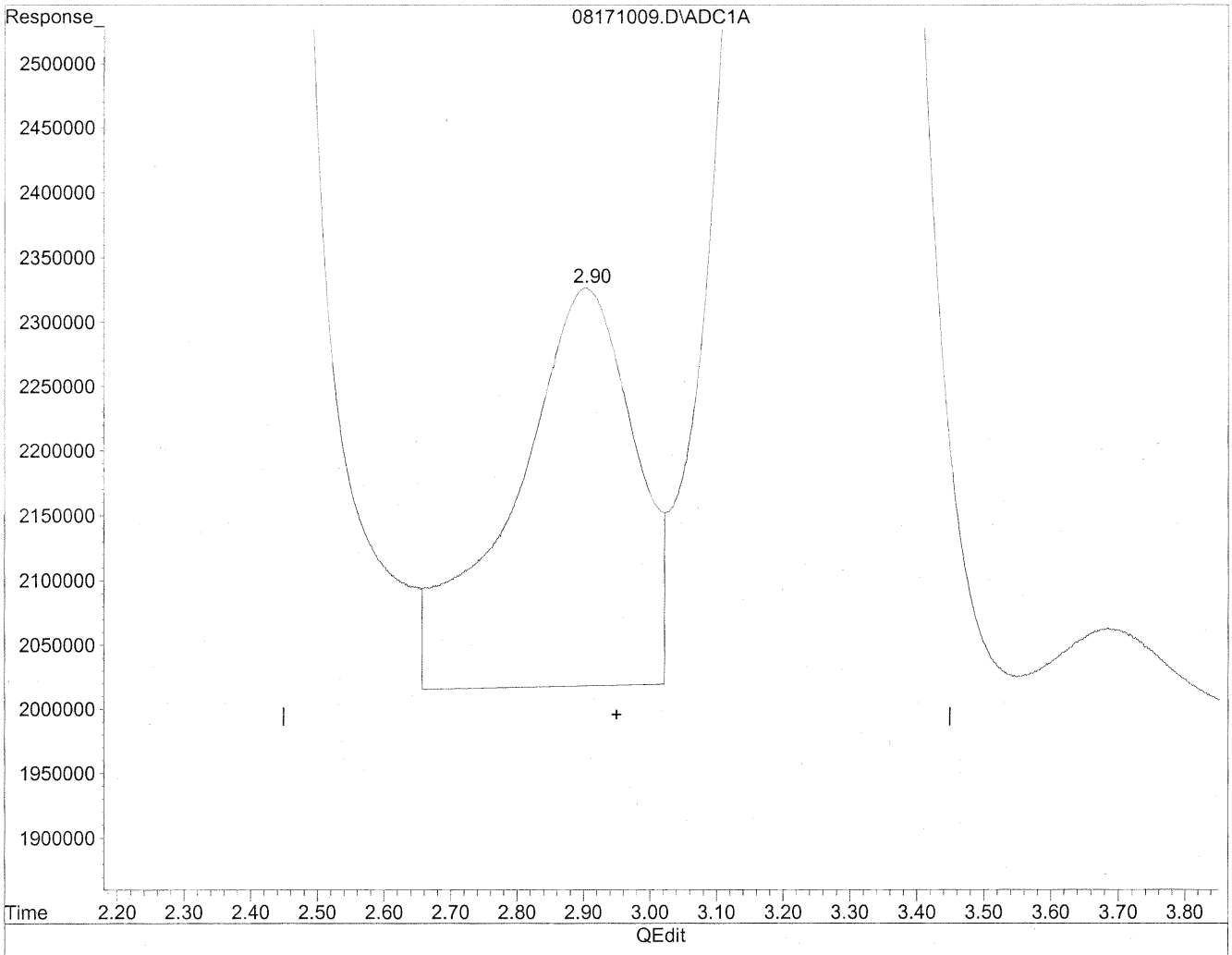
(+) = Expected Retention Time

08171009.D TO110709.M Sat Aug 22 14:18:53 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.90min 365.385ng/ml

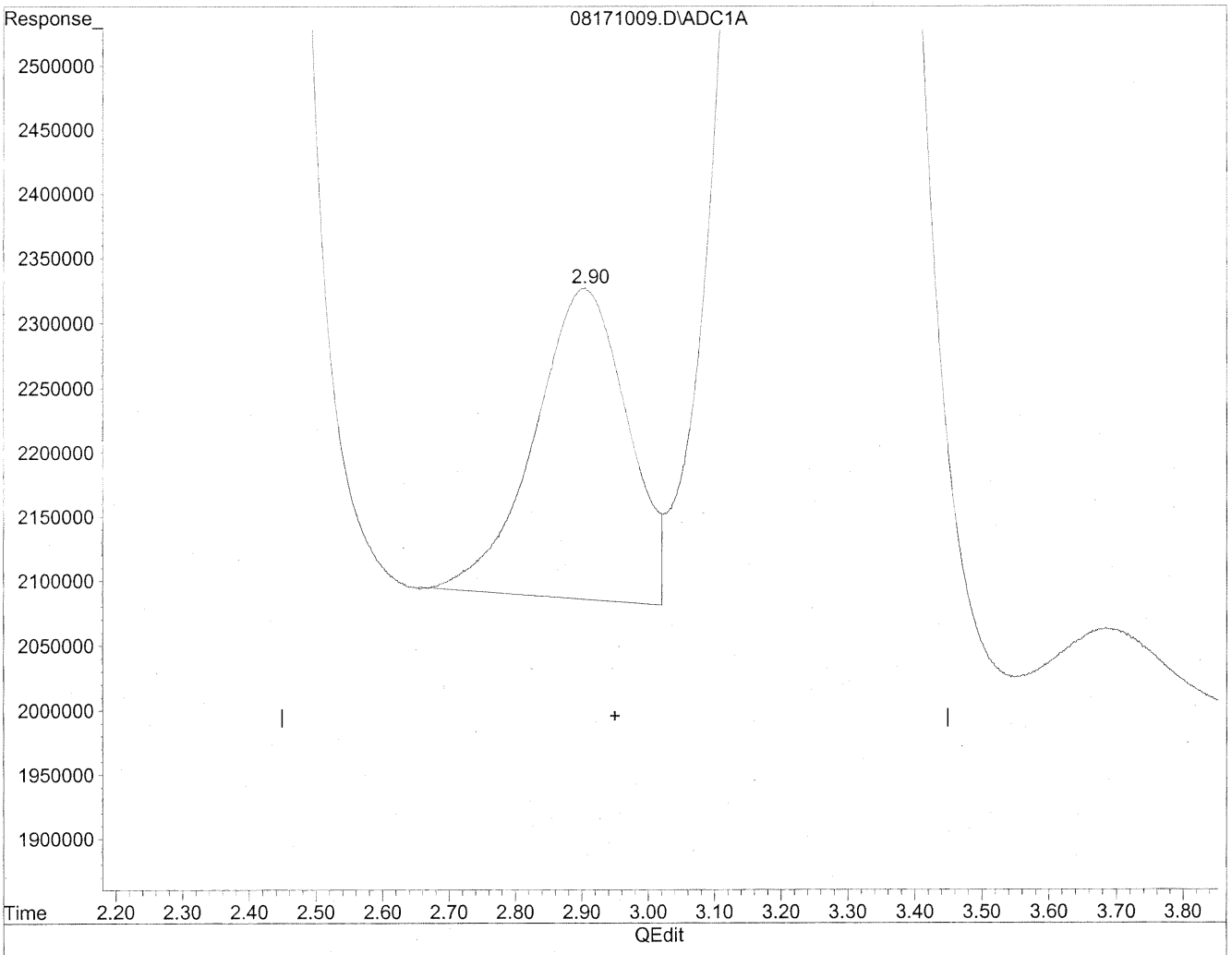
response 38984860

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.90min 220.236ng/ml m

response 23498130

HC 8/22/09
BC

KP 8/23/09

90

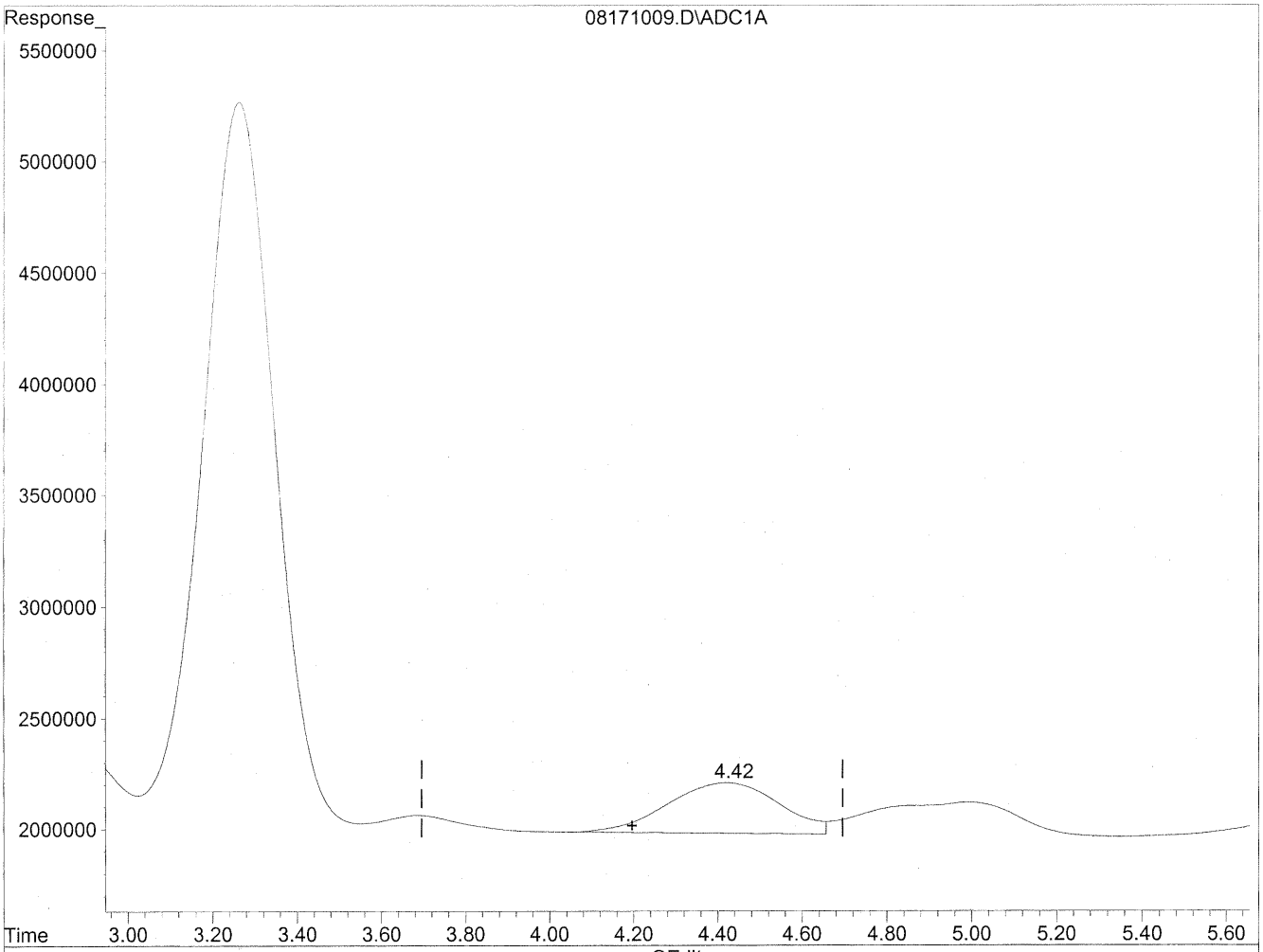
(+) = Expected Retention Time

08171009.D TO110709.M Sat Aug 22 14:19:12 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.42min 429.029ng/ml

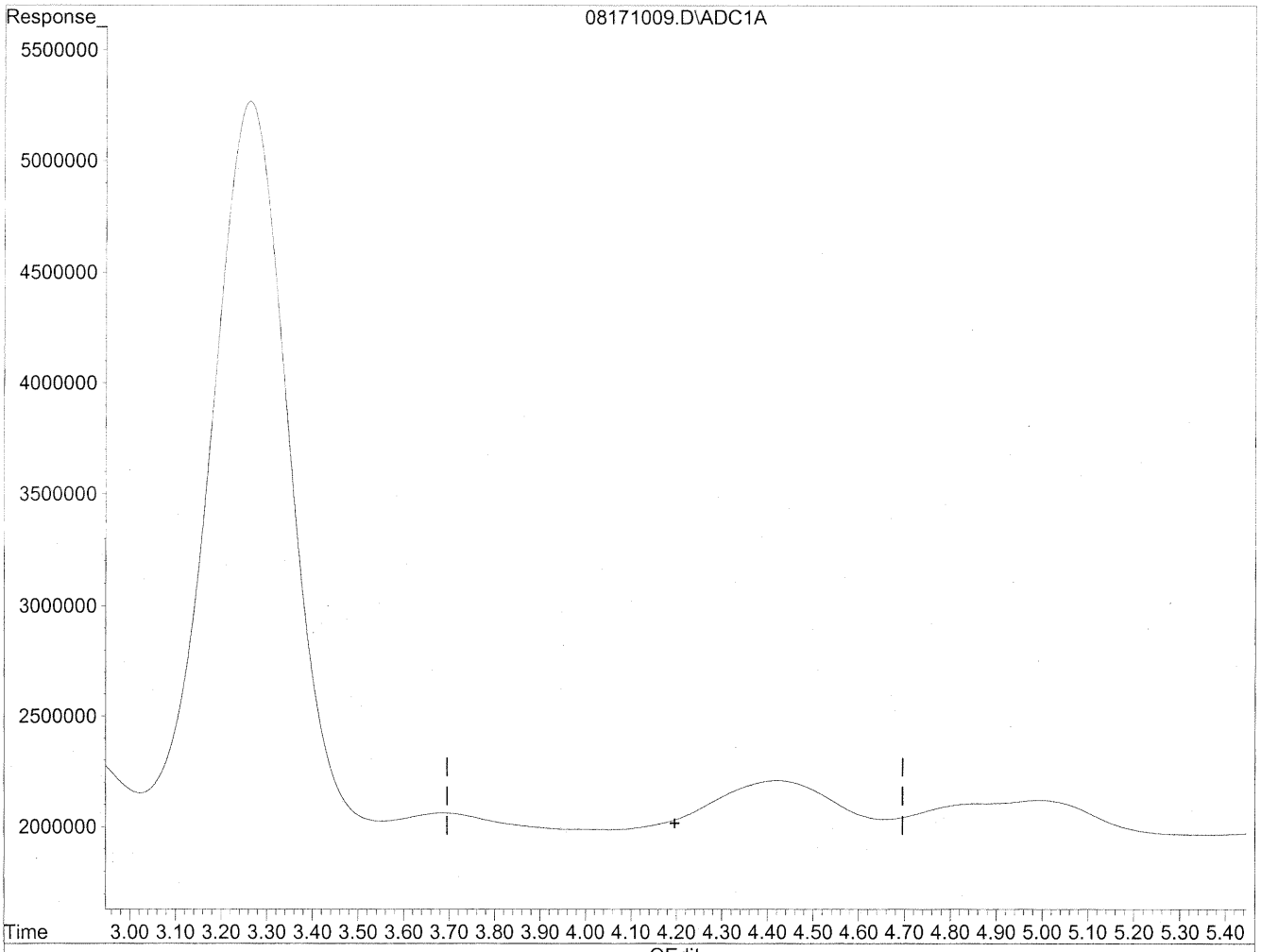
response 41793983

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

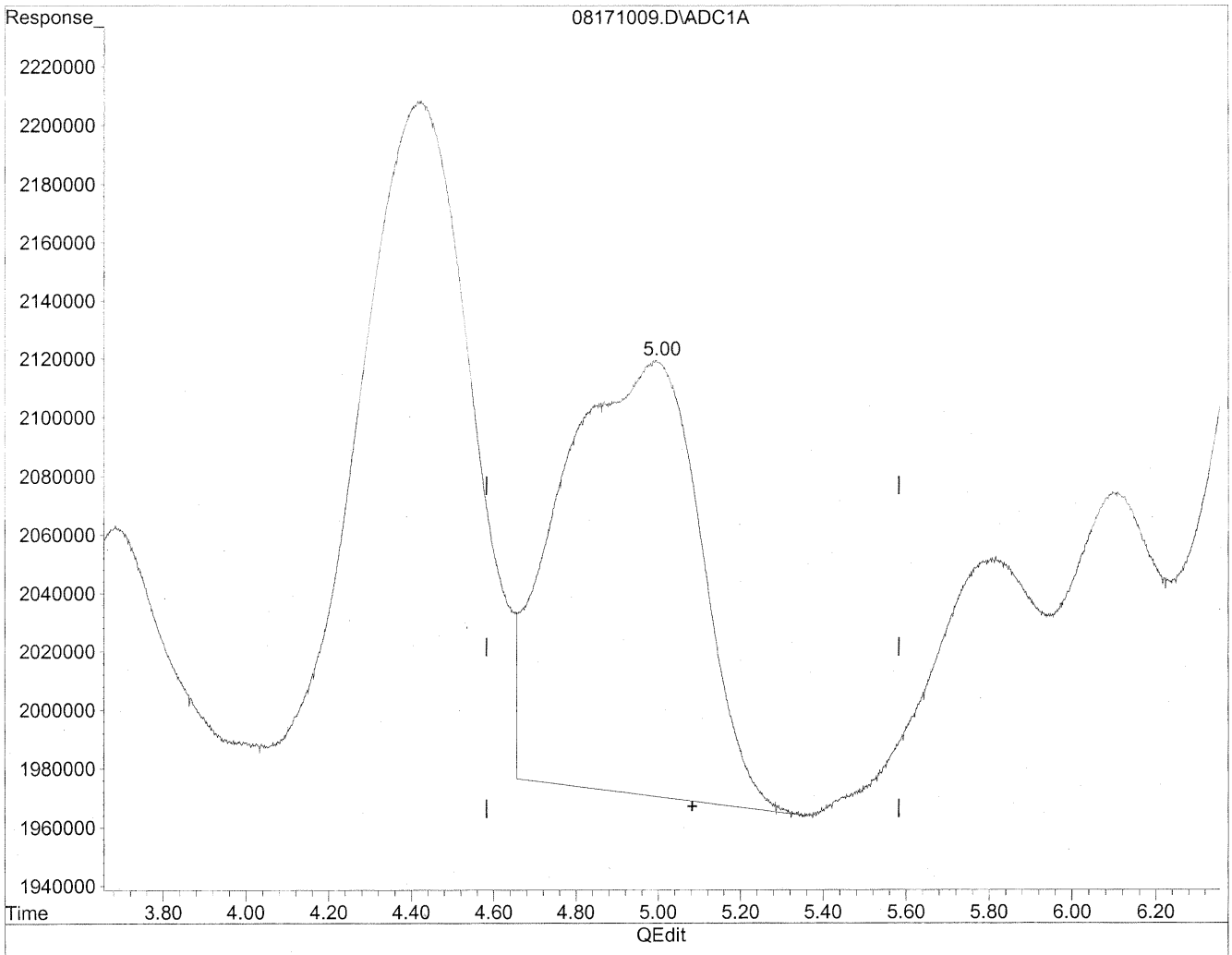
HC 8/22/09
MP
KR 8/23/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.99min 394.416ng/ml

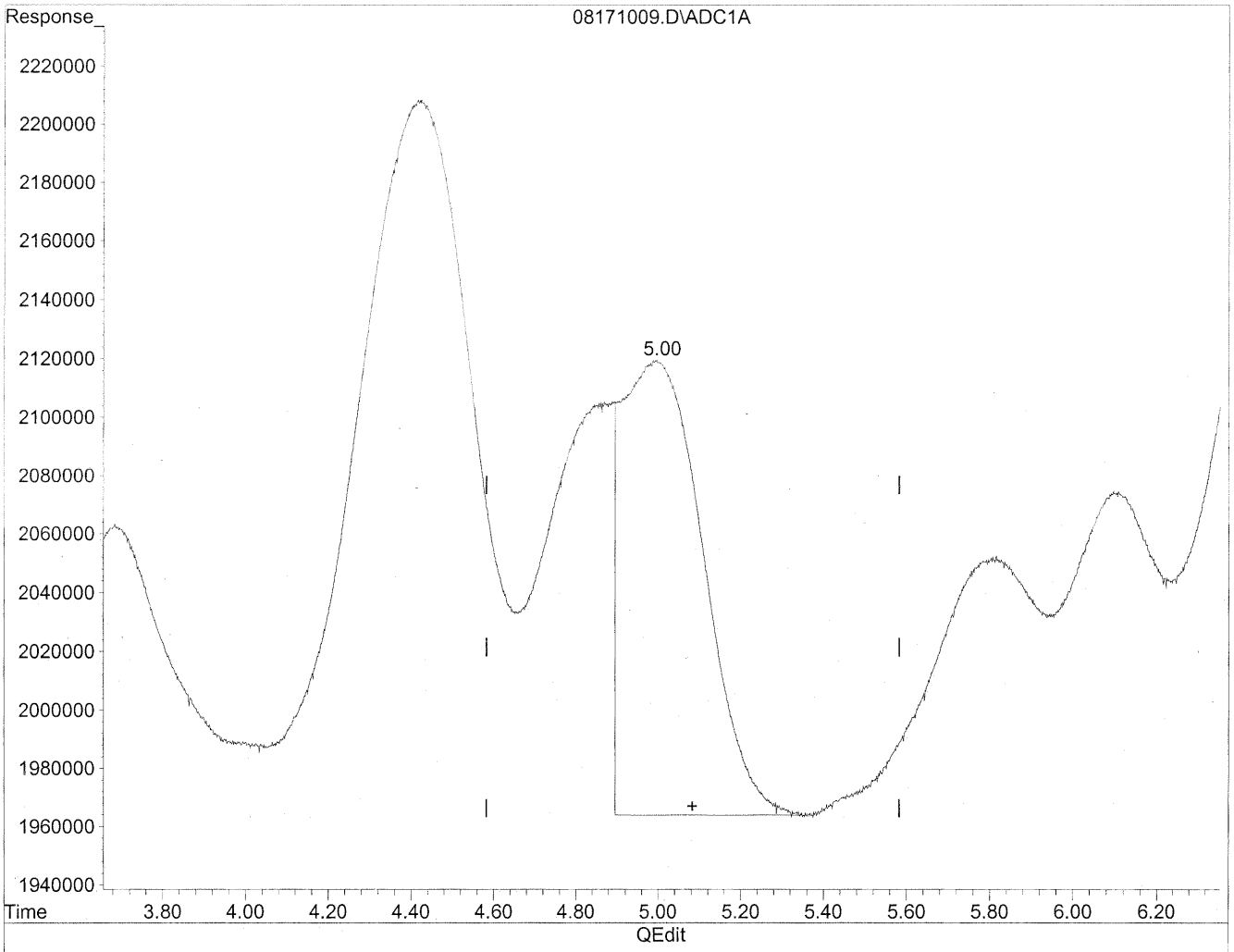
response 34841185

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.00min 240.626ng/ml m

response 21256004

tc 8/22/09
SP

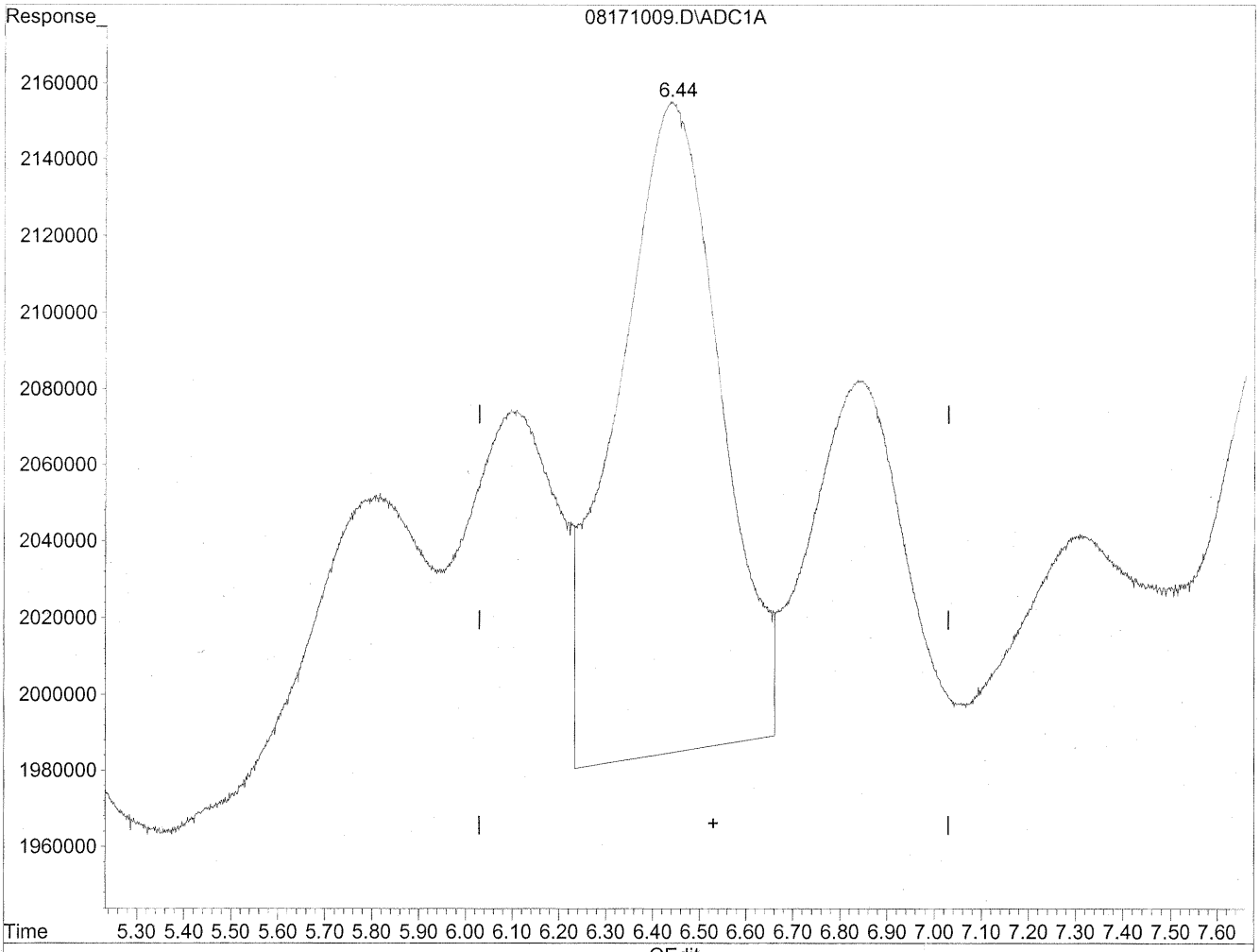
KP 8/23/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.44min 396.943ng/ml

response 26146346

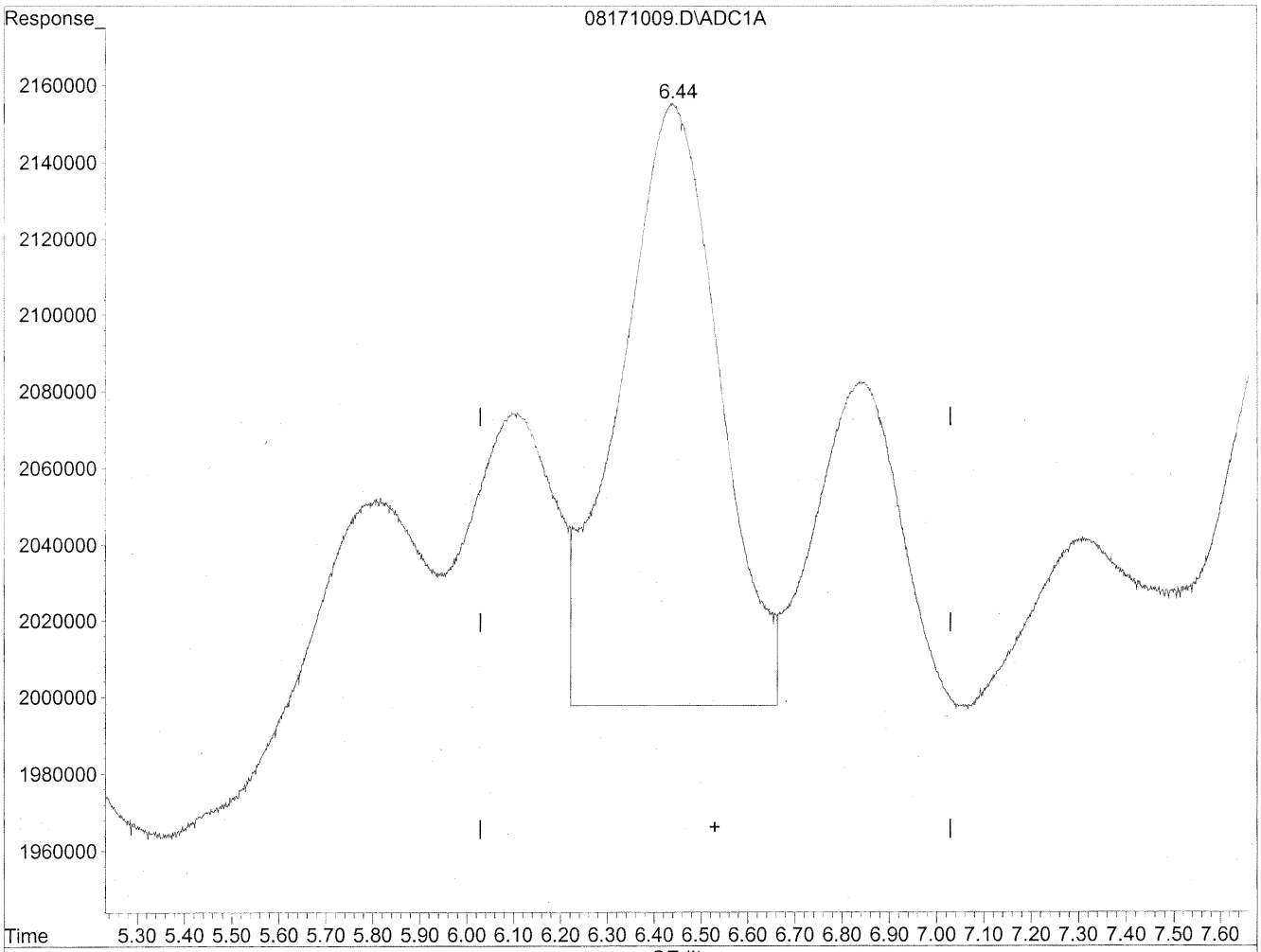
(+) = Expected Retention Time

08171009.D TO110709.M Sat Aug 22 14:19:43 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.44min 352.343ng/ml m

response 23208609

HC 8/22/09
BC
8/23/09

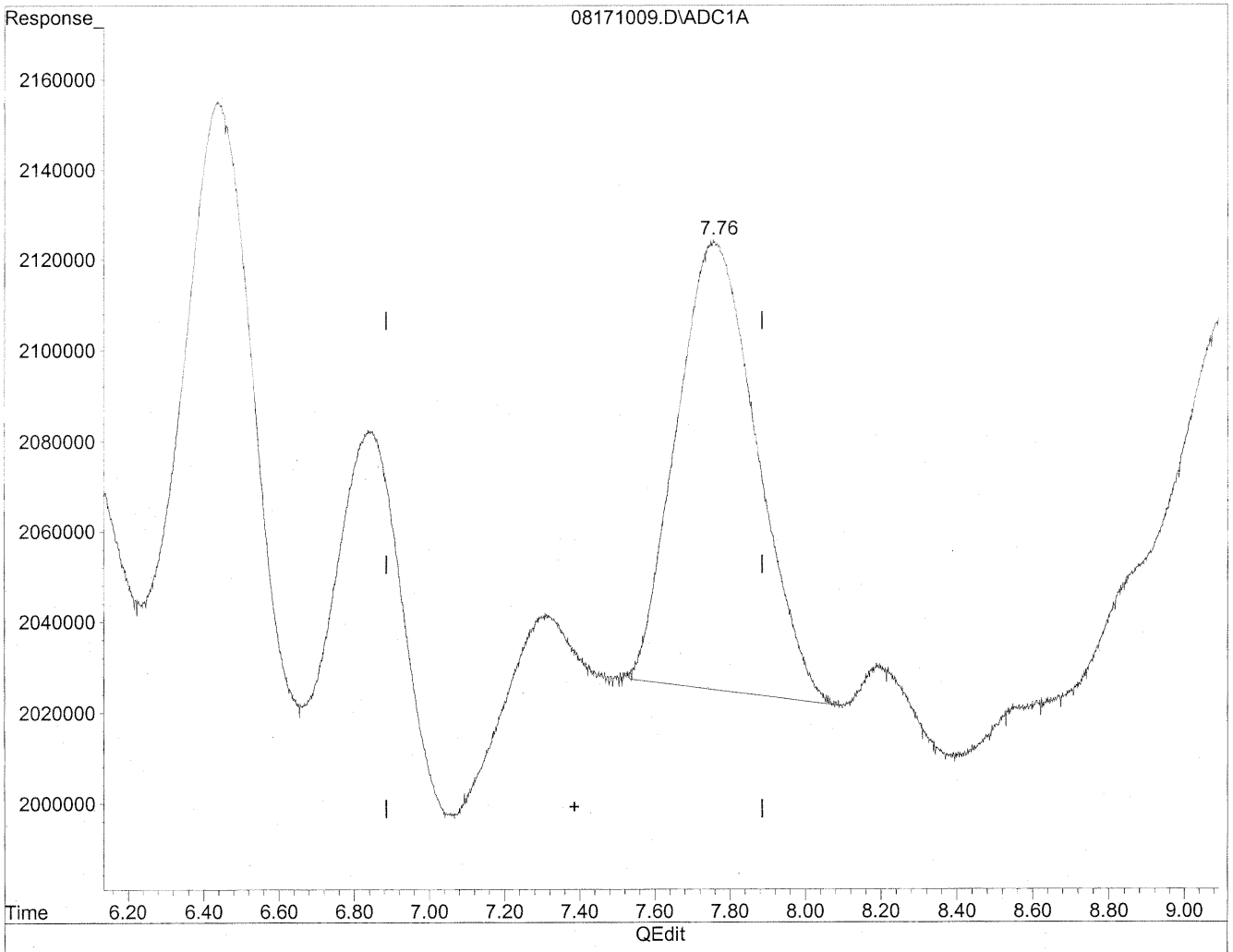
(+) = Expected Retention Time

08171009.D TO110709.M Sat Aug 22 14:20:11 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.76min 187.013ng/ml

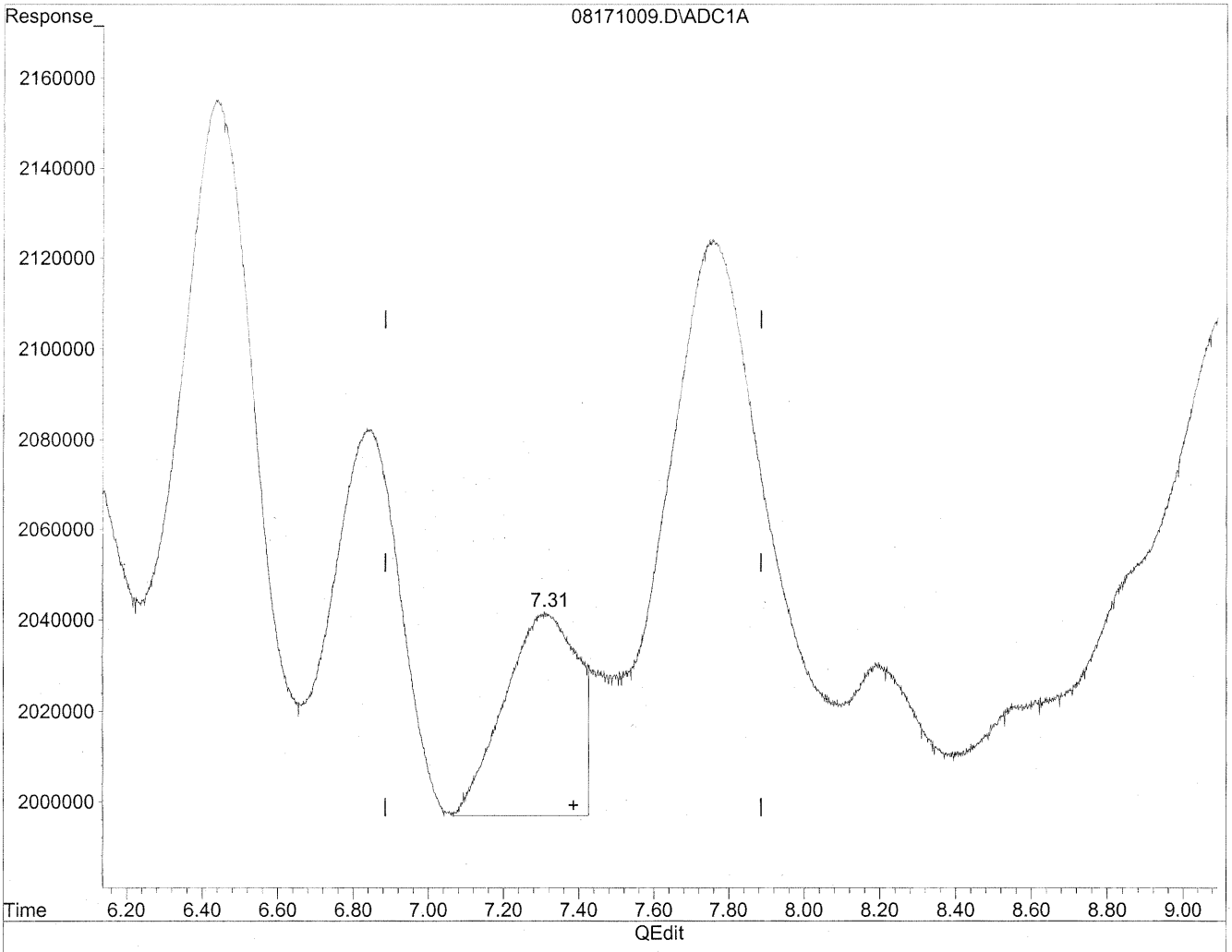
response 14633955

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.31min 76.853ng/ml m

response 6013814

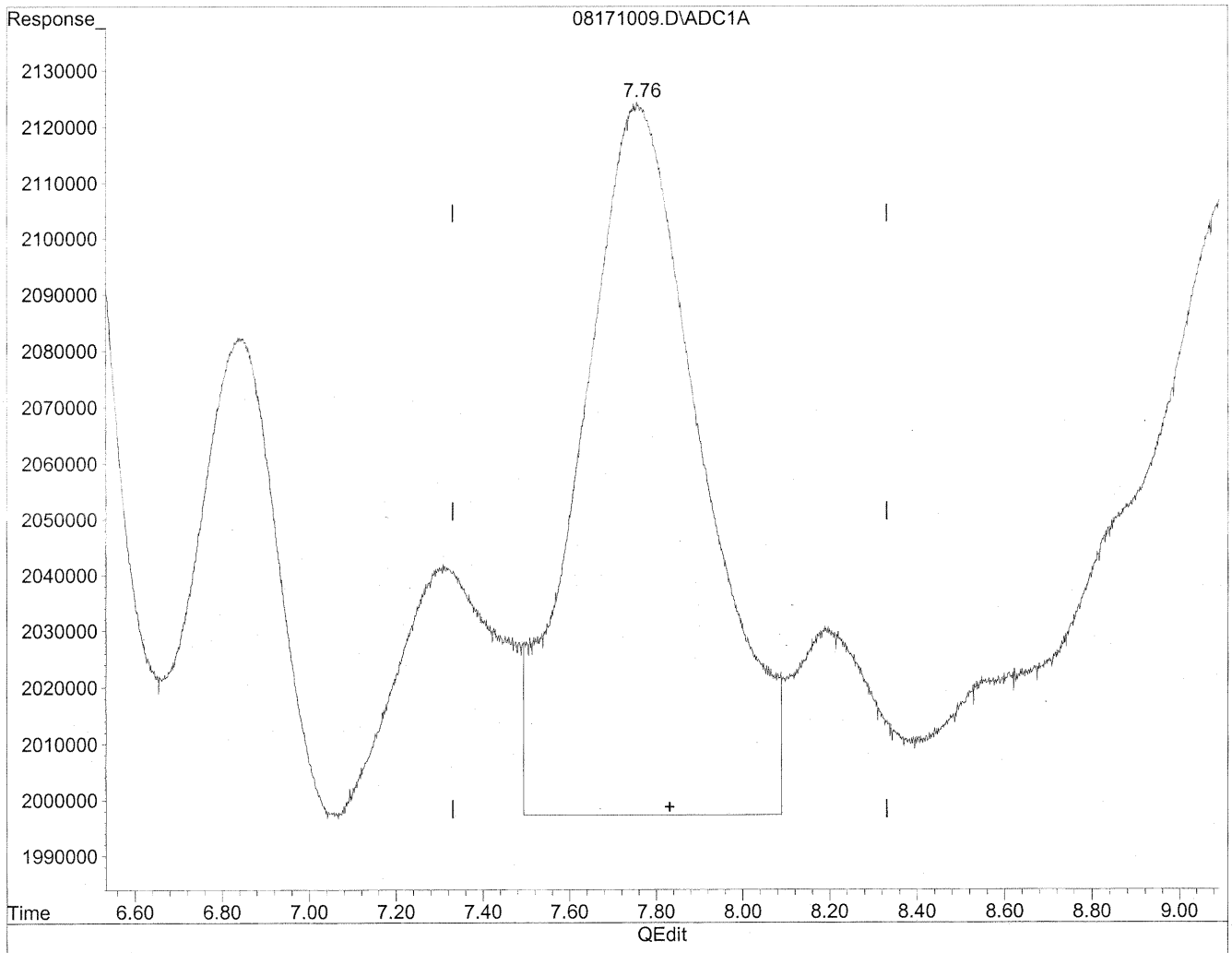
HC 8/22/09 MP

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.76min 332.014ng/ml m

response 24404666

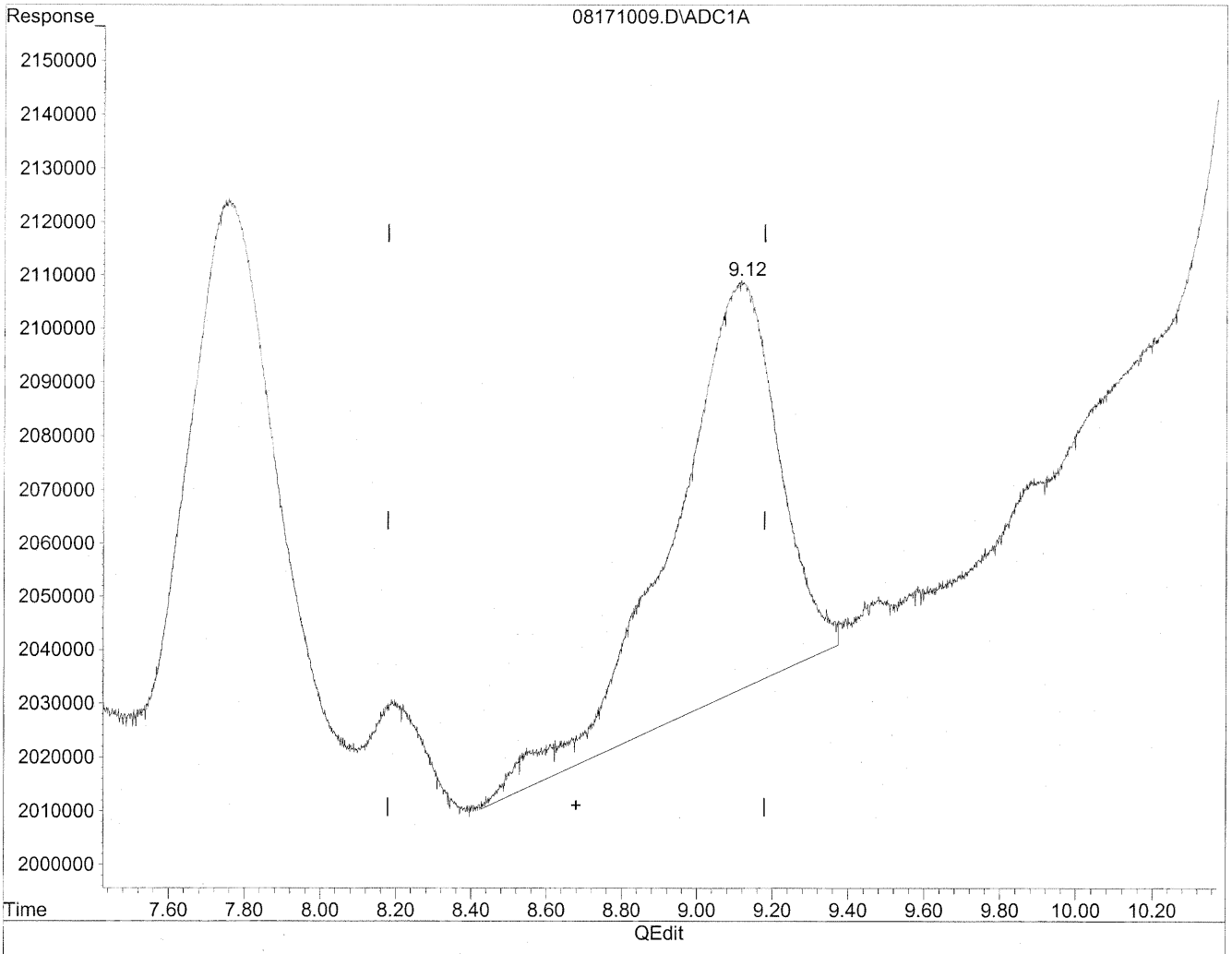
HC 8/22/09
1 c
no release
KP 8/25/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.12min 255.926ng/ml

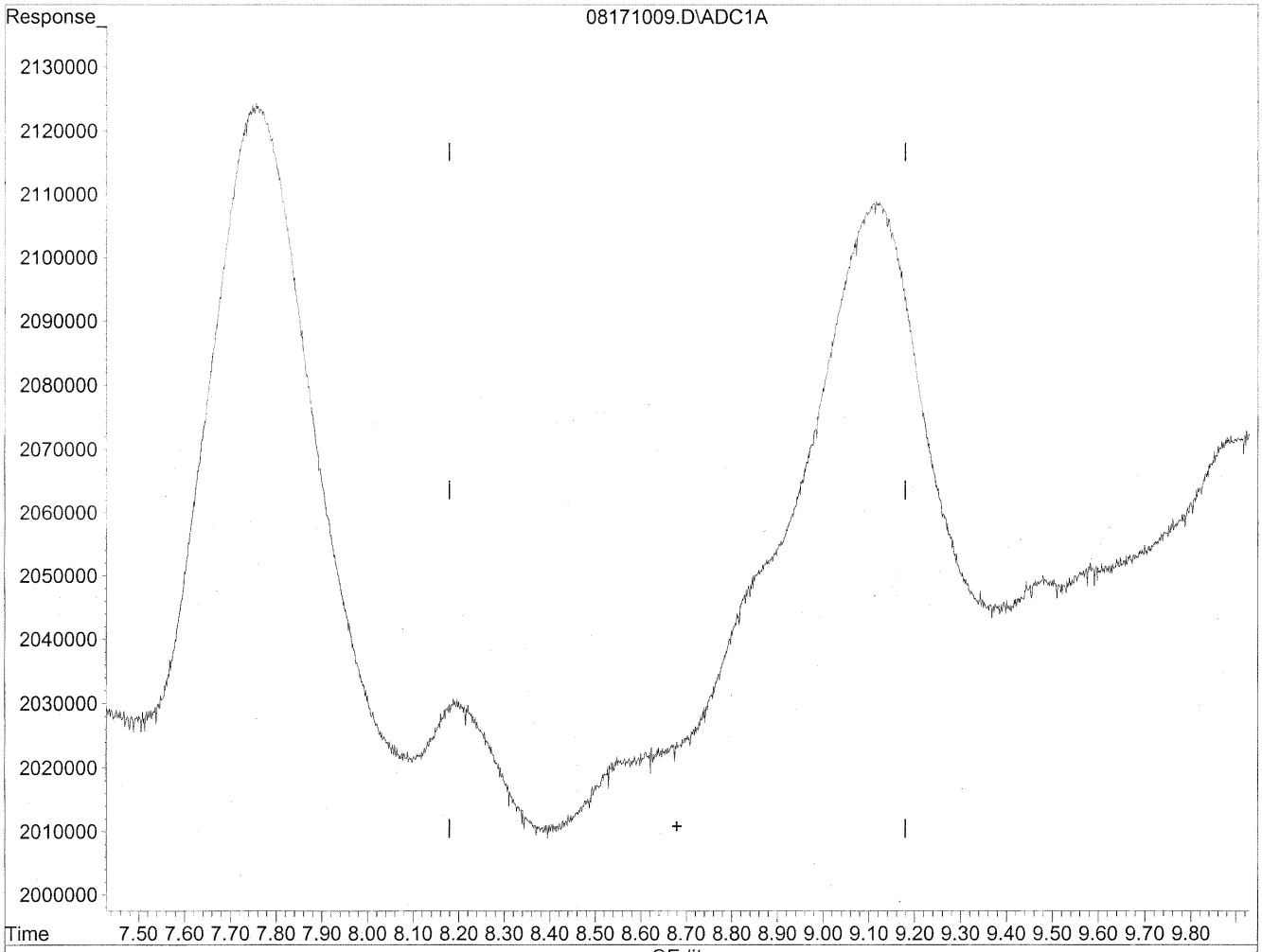
response 14925694

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09 MP

HC 8/23/09

101

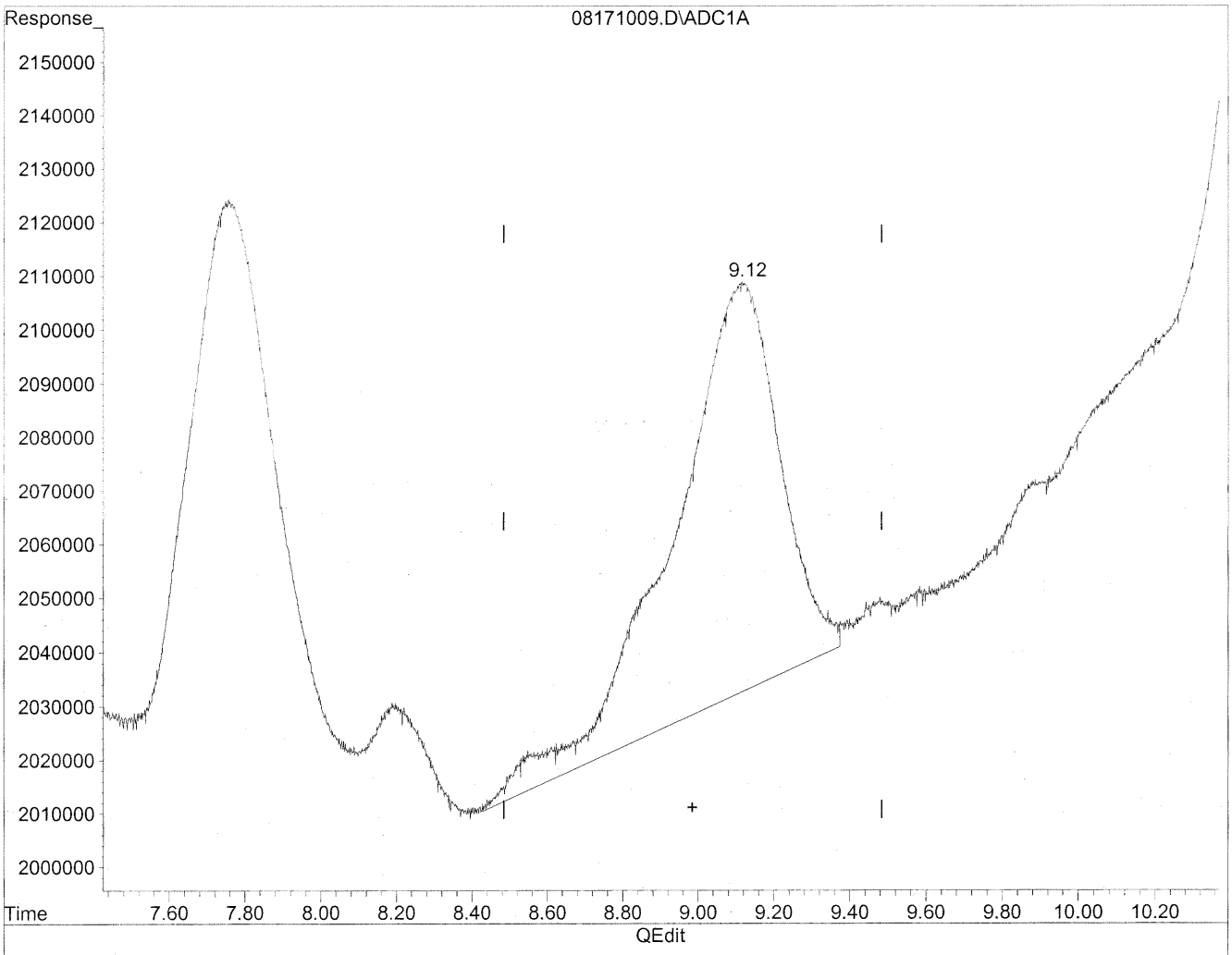
(+) = Expected Retention Time

08171009.D TO110709.M Sat Aug 22 14:20:54 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

9.12min 276.425ng/ml

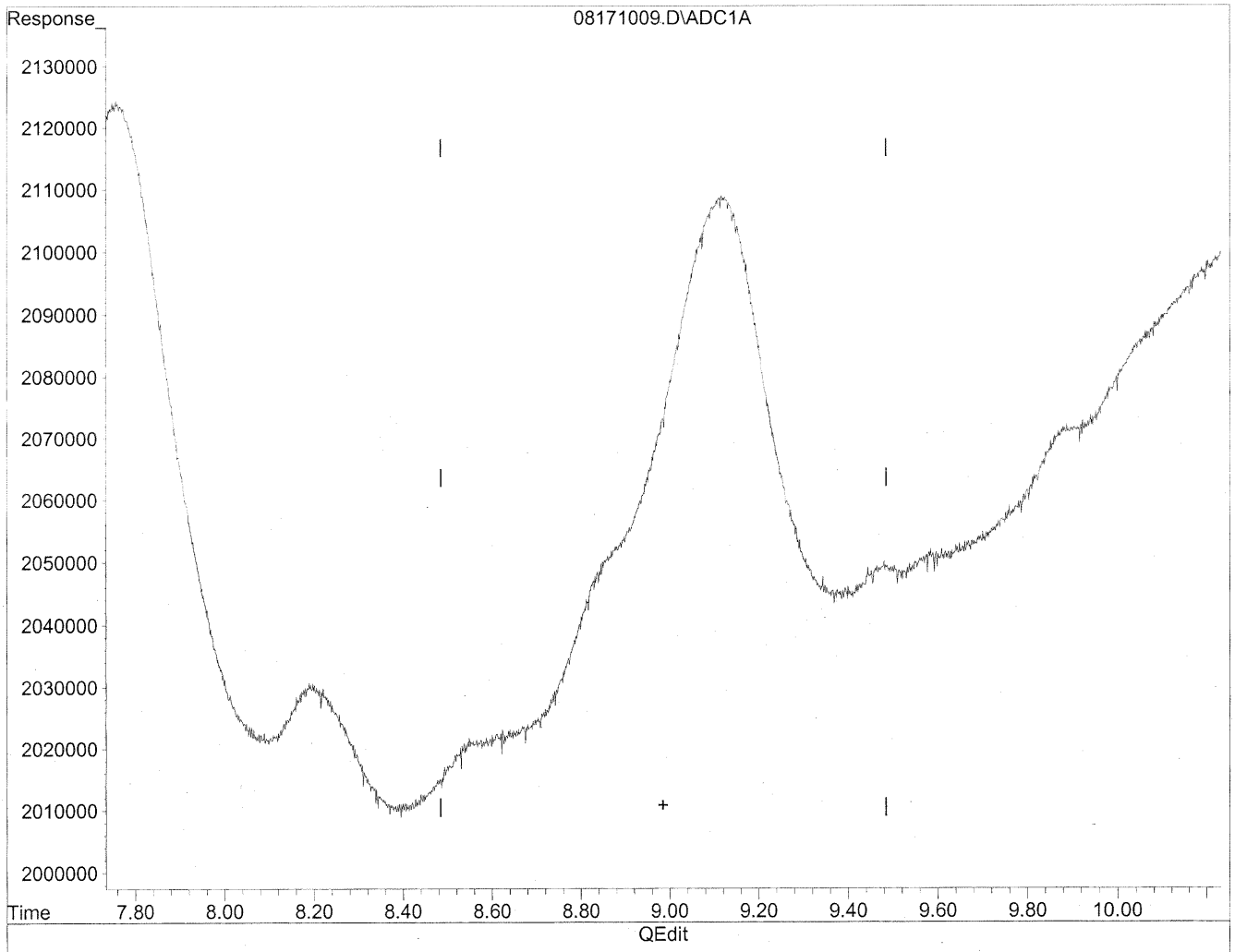
response 14925694

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

0.00min 0.000ng/ml d

response 0

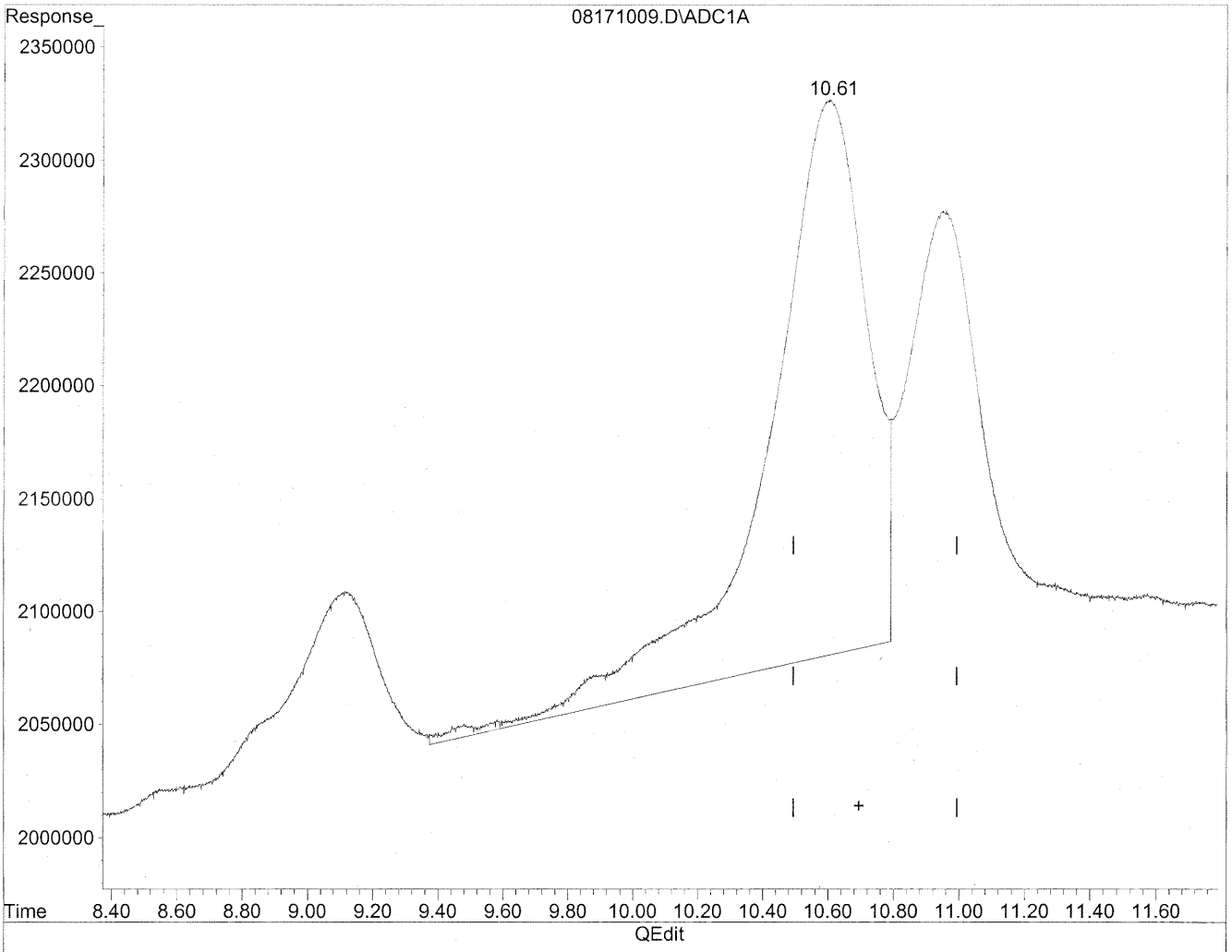
HC 8/22/09
MP
KP 8/23/09

(+) = Expected Retention Time
08171009.D TO110709.M Sat Aug 22 14:21:01 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.60min 766.131ng/ml

response 51594162

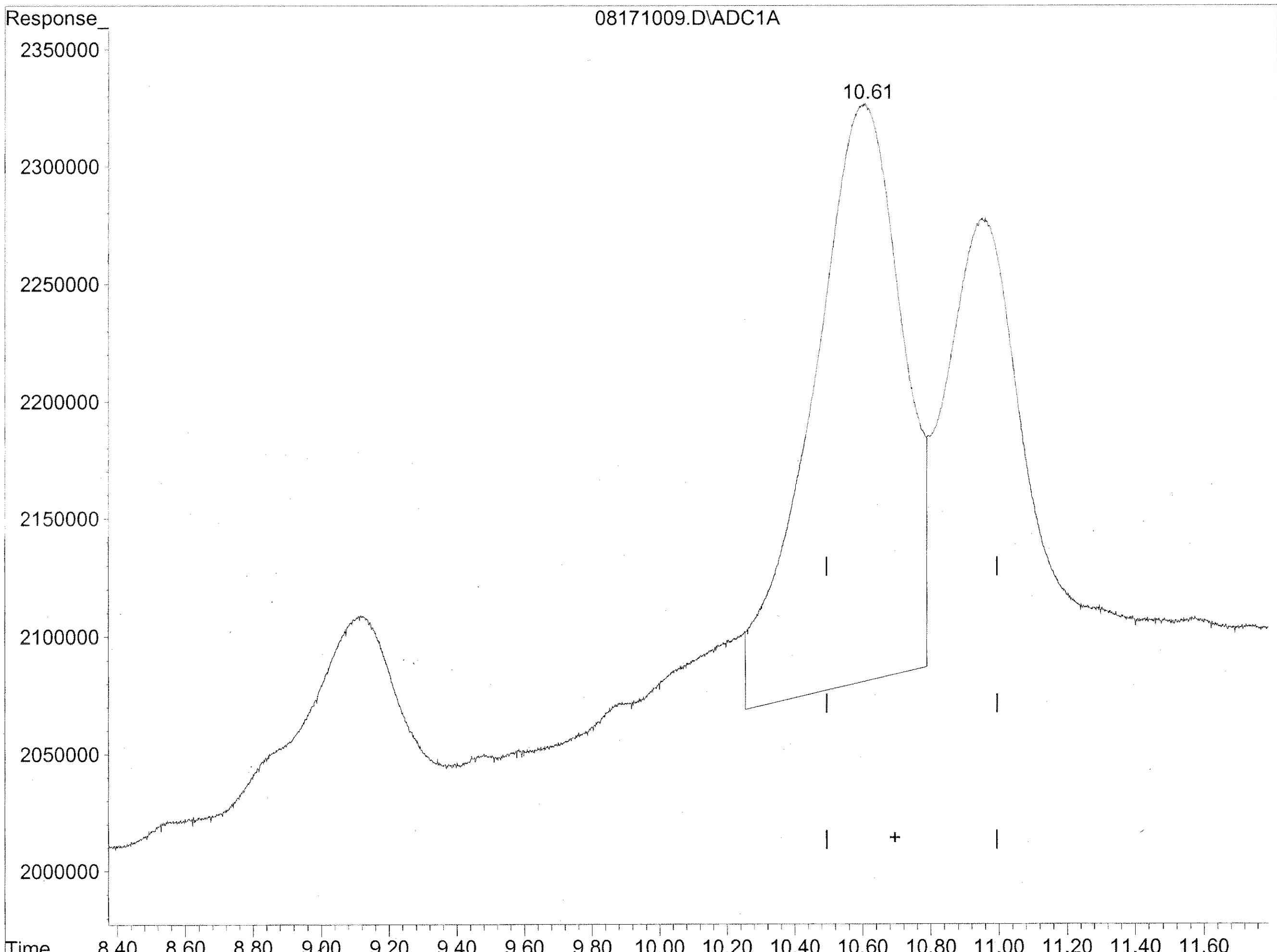
(+) = Expected Retention Time

08171009.D TO110709.M Sat Aug 22 14:21:05 2009

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D      Vial: 22
Acq On    : 18 Aug 2009   5:54 pm                    Operator: HC
Sample    : P0902772-004 front 1.0ml                 Inst    : LC 01
Misc      :                                          Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Aug 19  9:19 19109  Quant Results File: TO110709.RES

Method         : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title          : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update    : Wed Aug 19 10:45:48 2009
Response via   : Multiple Level Calibration
```



QEdit

(11) Hexaldehyde

10.61min 669.492ng/ml m

response 45086096

HC 8/21/09 ✓

KR 8/23/09

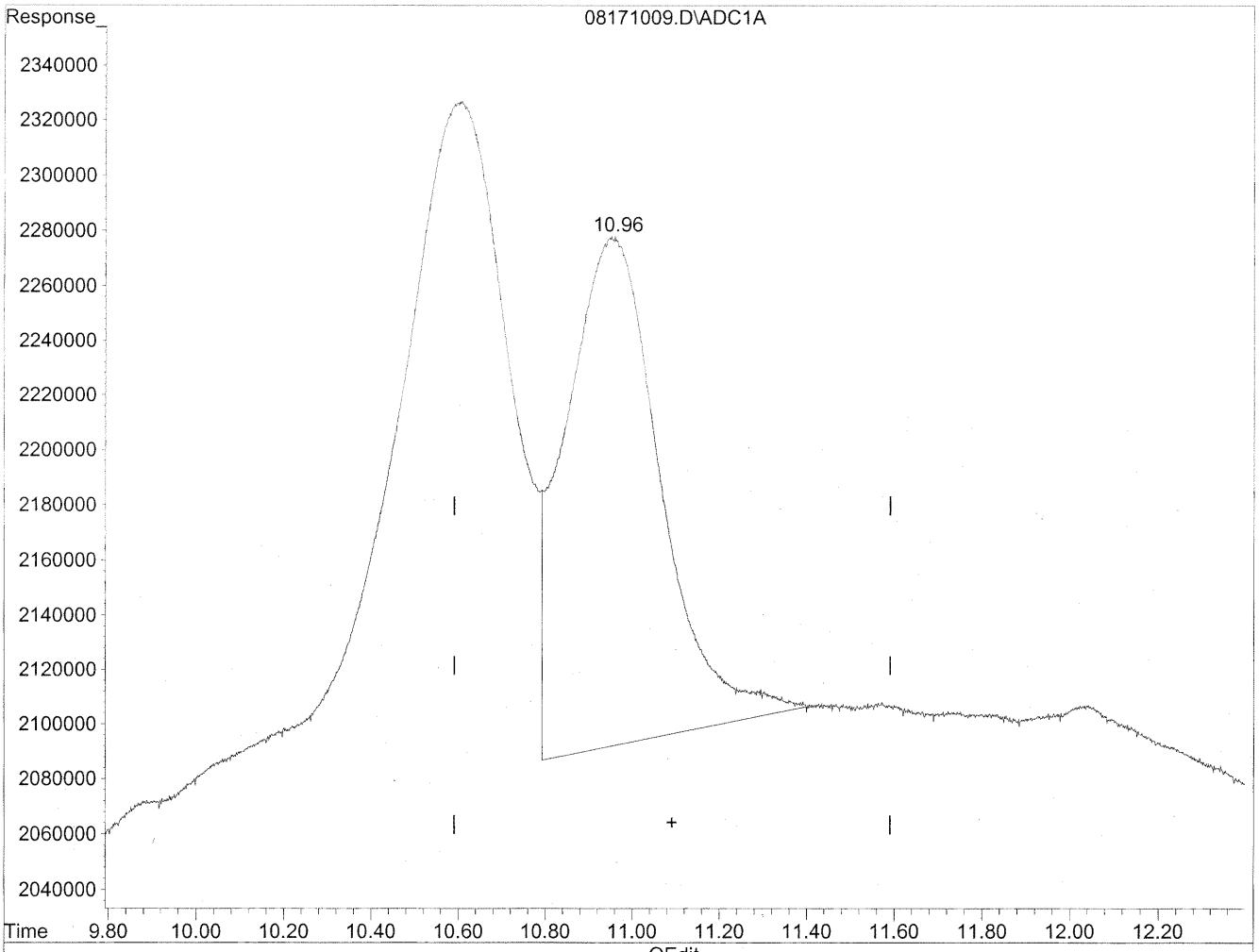
(+) = Expected Retention Time

08171009.D TO110709.M Sat Aug 22 14:21:14 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171009.D Vial: 22
Acq On : 18 Aug 2009 5:54 pm Operator: HC
Sample : P0902772-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.95min 577.570ng/ml

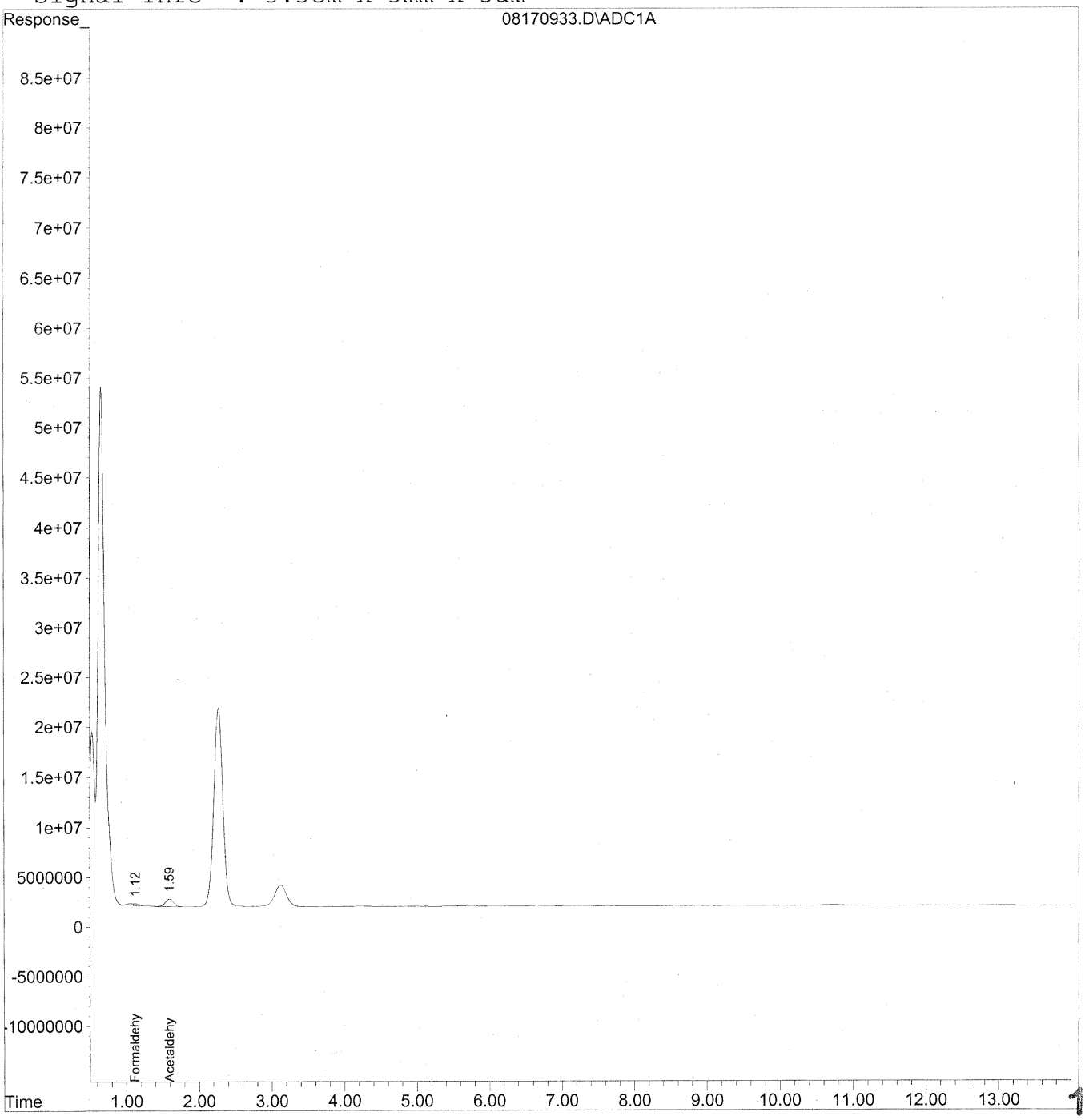
response 28308687

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_08\17\08170933.D Vial: 32
Acq On : 17 Aug 2009 10:51 pm Operator: HC
Sample : P0902772-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:14 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



107

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170933.D Vial: 32
Acq On : 17 Aug 2009 10:51 pm Operator: HC
Sample : P0902772-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:14 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

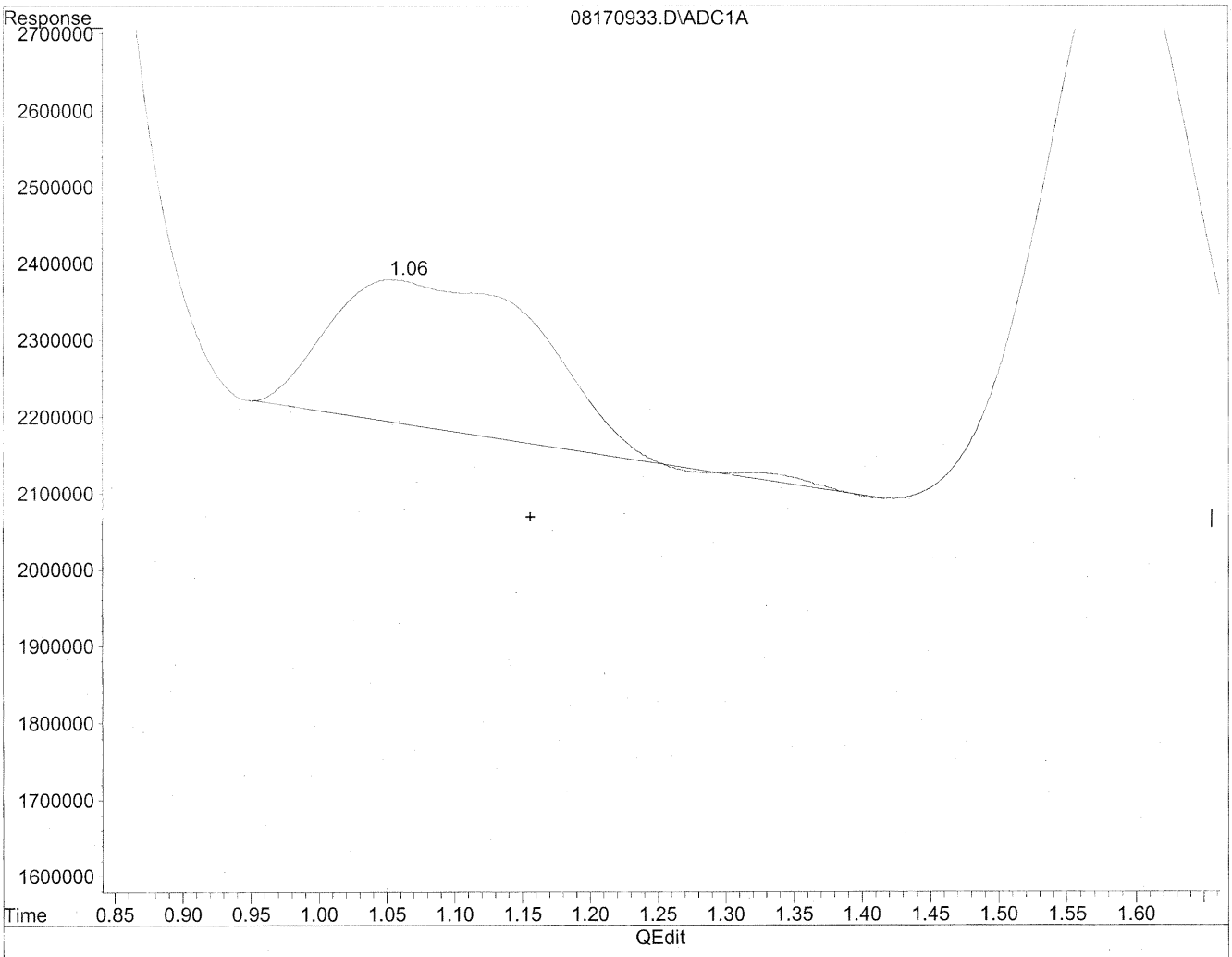
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.12	10660553	58.070	ng/mlm
2)	Acetaldehyde	1.59	59681352	425.616	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/mld
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/mld

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170933.D Vial: 32
Acq On : 17 Aug 2009 10:51 pm Operator: HC
Sample : P0902772-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.05min 115.972ng/ml

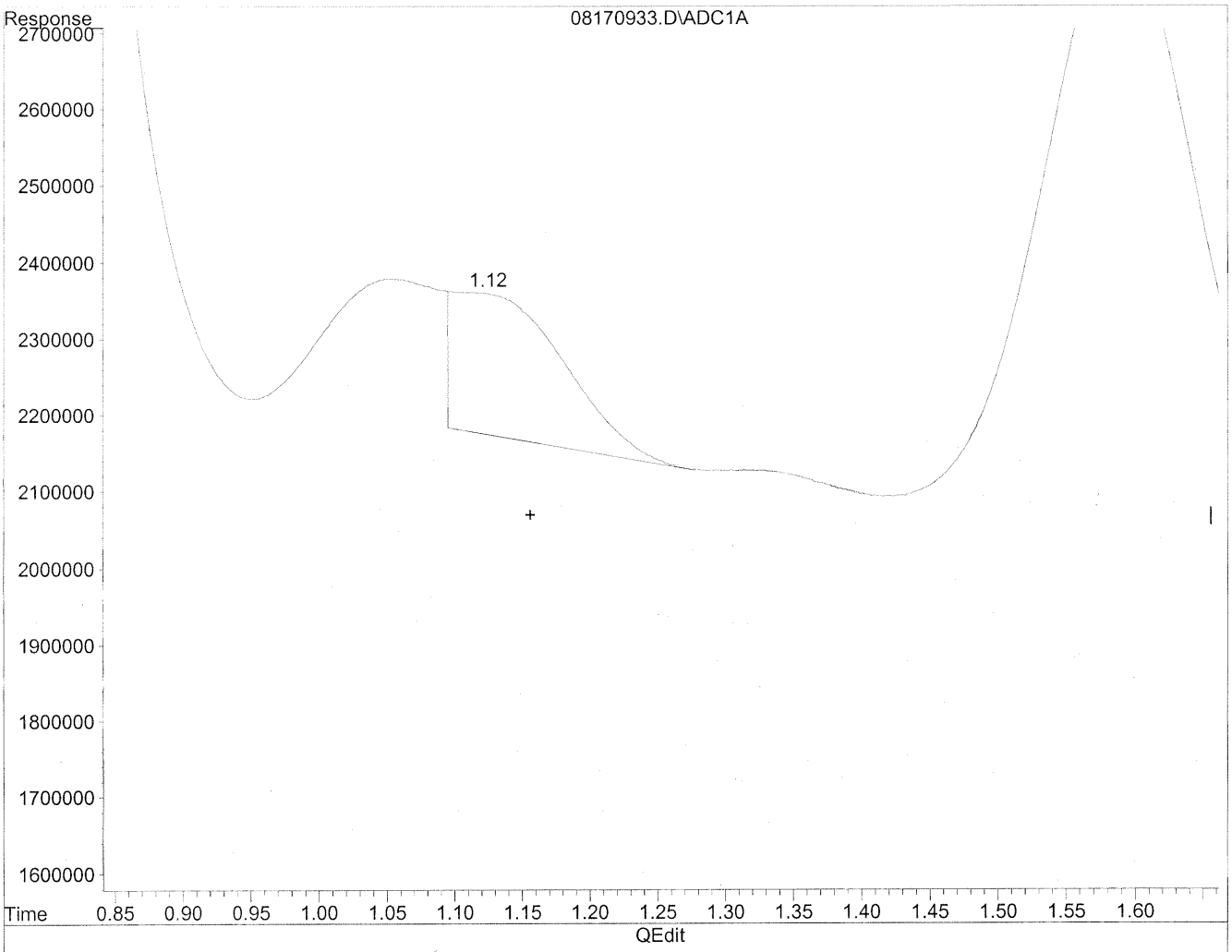
response 21290391

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170933.D Vial: 32
Acq On : 17 Aug 2009 10:51 pm Operator: HC
Sample : P0902772-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.12min 58.070ng/ml m

response 10660553

HC 8/21/09
SP

KRB 8/23/09

110

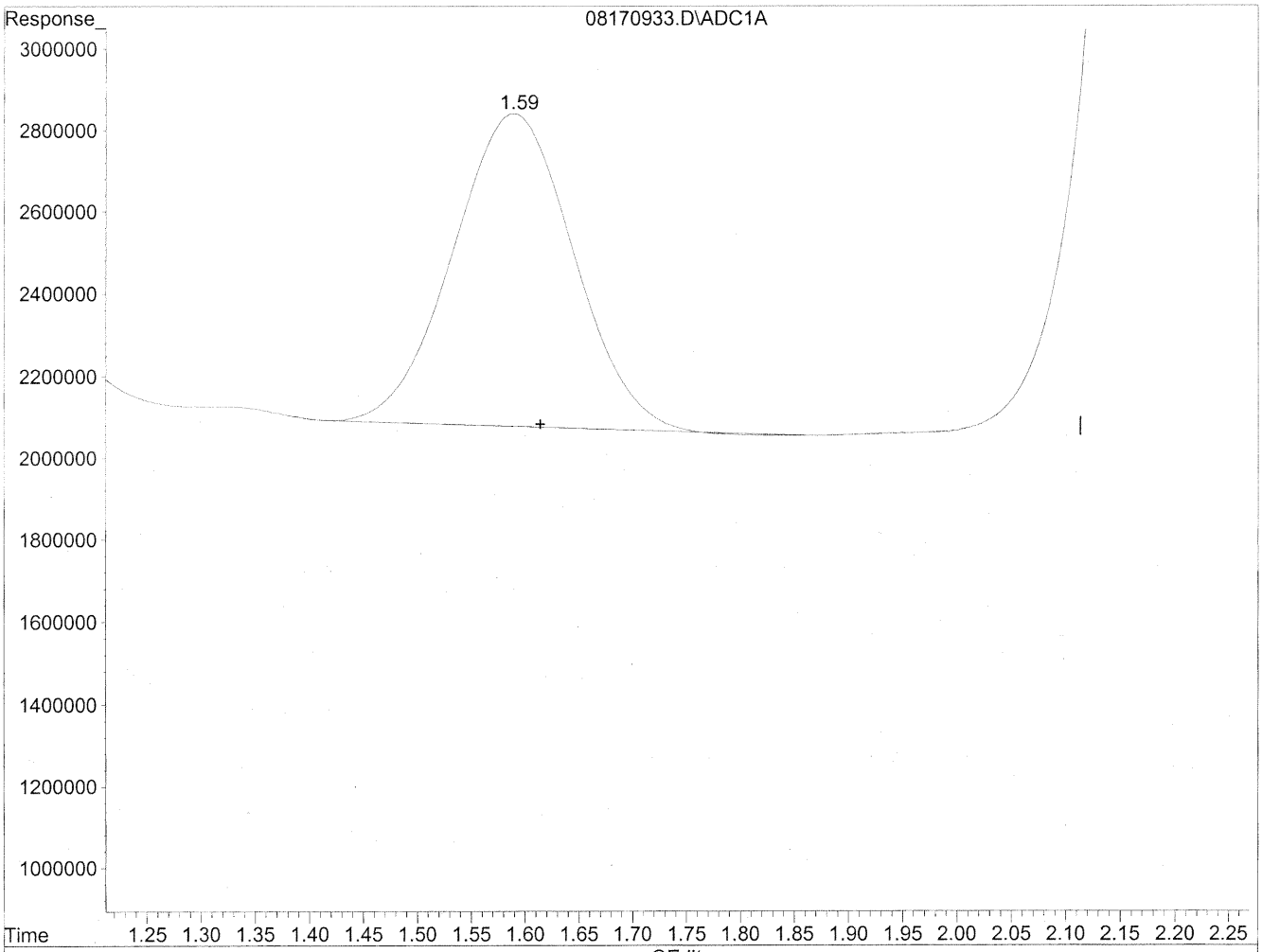
(+) = Expected Retention Time

08170933.D TO110709.M Fri Aug 21 18:14:03 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170933.D Vial: 32
Acq On : 17 Aug 2009 10:51 pm Operator: HC
Sample : P0902772-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.59min 422.071ng/ml

response 59184227

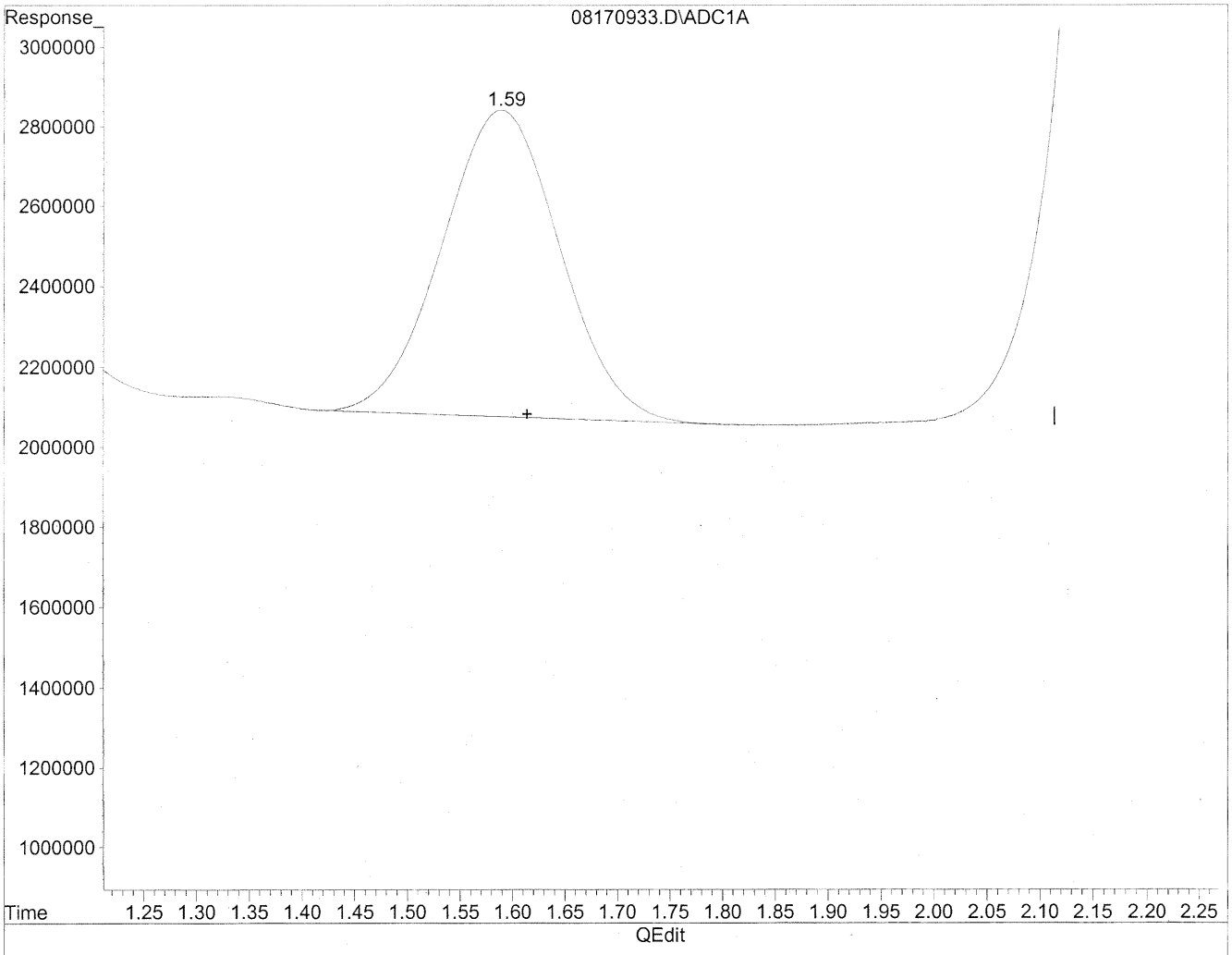
111

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170933.D Vial: 32
Acq On : 17 Aug 2009 10:51 pm Operator: HC
Sample : P0902772-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.59min 425.616ng/ml m

response 59681352

HC 8/24/09
LC

KR 8/23/09

112

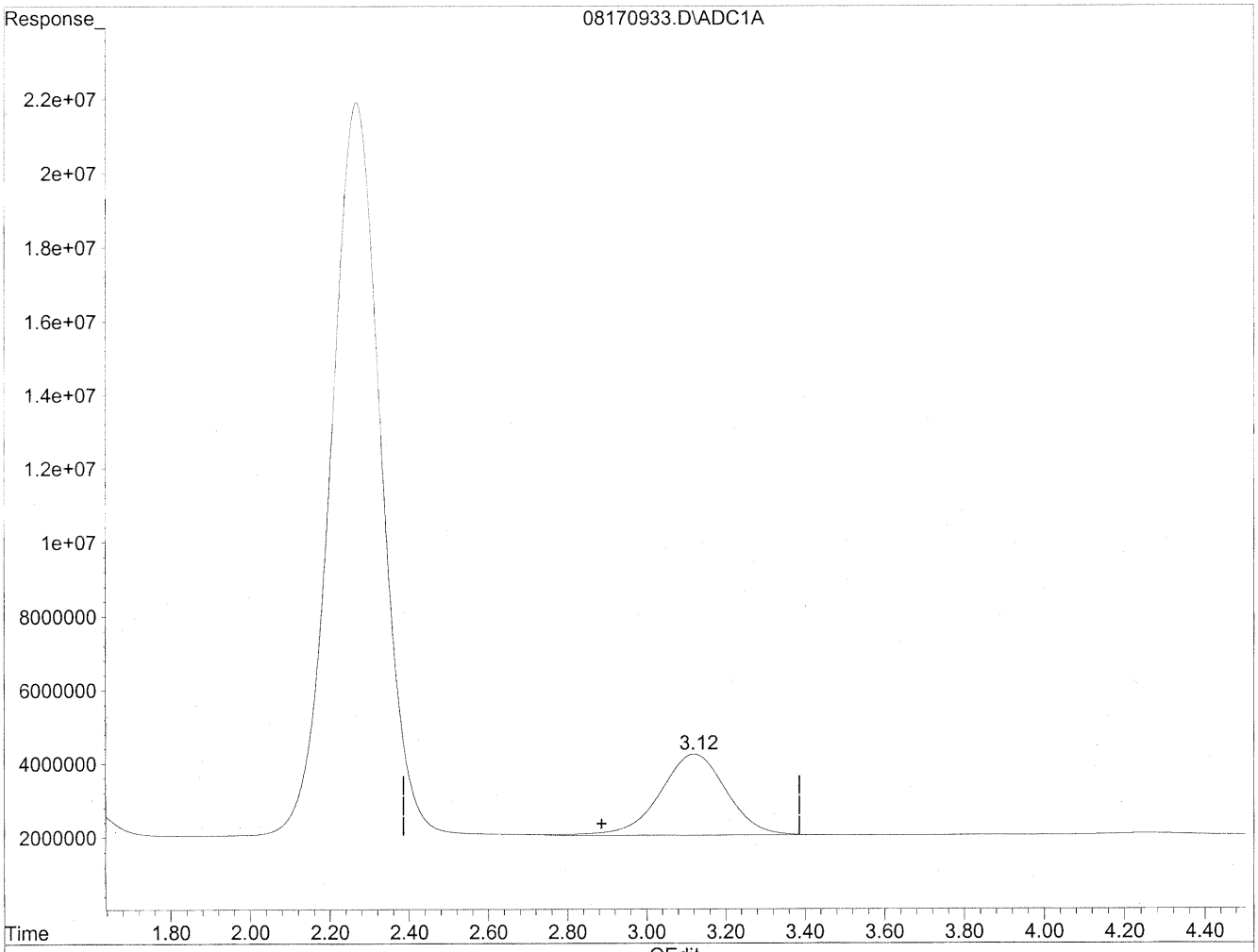
(+) = Expected Retention Time

08170933.D TO110709.M Fri Aug 21 18:14:26 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170933.D Vial: 32
Acq On : 17 Aug 2009 10:51 pm Operator: HC
Sample : P0902772-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.12min 2394.584ng/ml

response 255490717

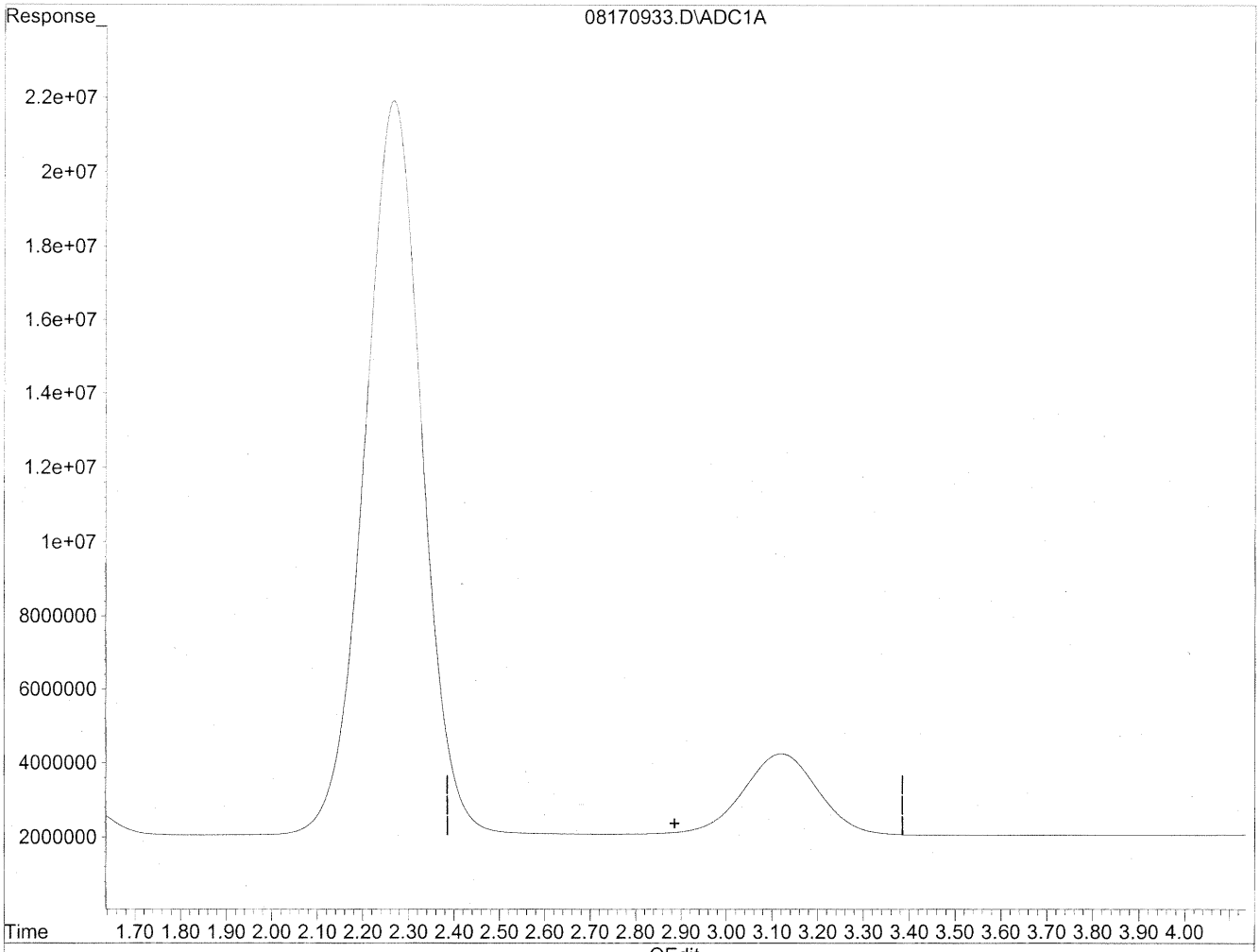
(+) = Expected Retention Time

08170933.D TO110709.M Fri Aug 21 18:14:29 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170933.D Vial: 32
Acq On : 17 Aug 2009 10:51 pm Operator: HC
Sample : P0902772-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HL 8/24/09 WP

KP 8/23/07

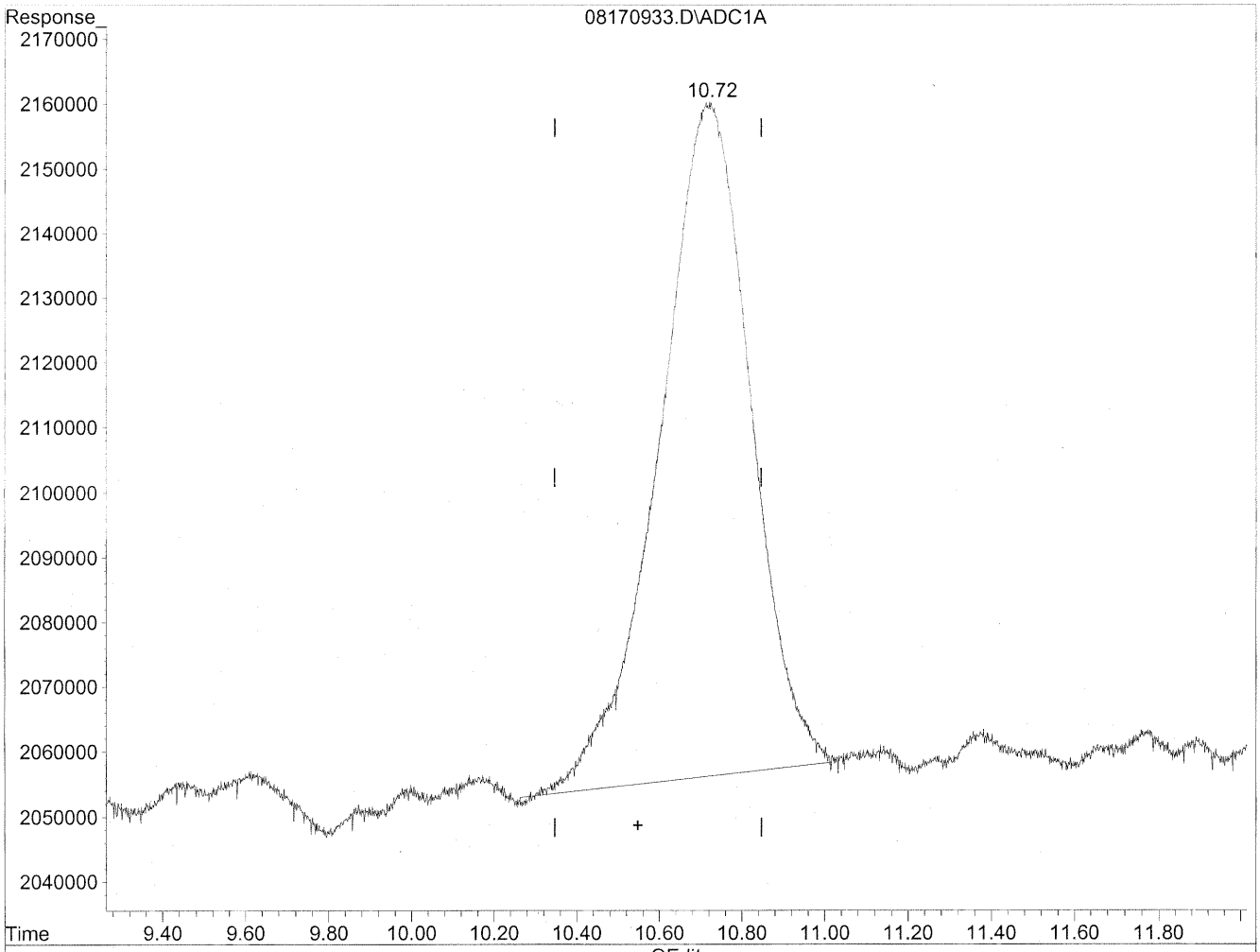
114

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170933.D Vial: 32
Acq On : 17 Aug 2009 10:51 pm Operator: HC
Sample : P0902772-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.72min 236.648ng/ml

response 15936741

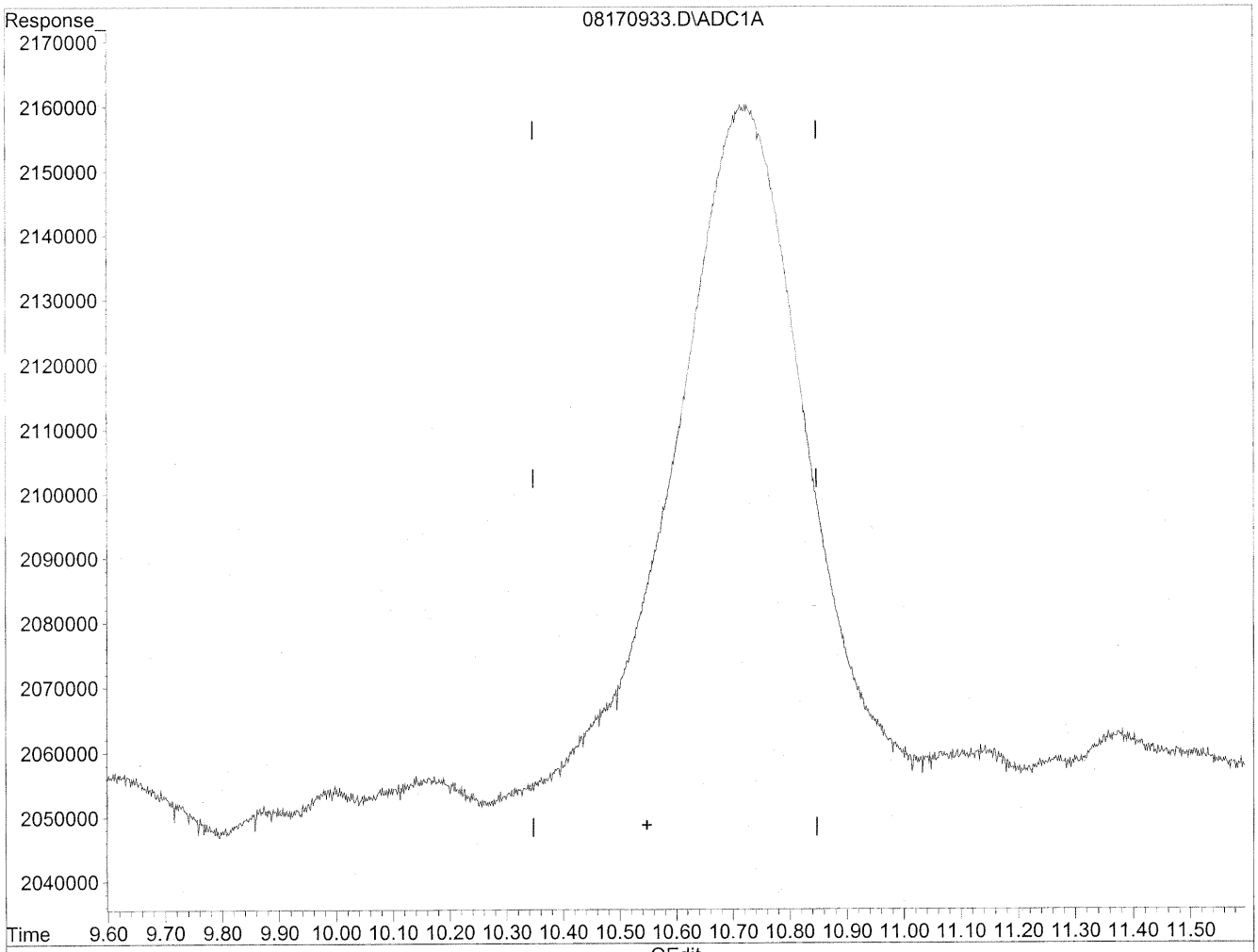
(+) = Expected Retention Time

08170933.D TO110709.M Fri Aug 21 18:14:37 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170933.D Vial: 32
Acq On : 17 Aug 2009 10:51 pm Operator: HC
Sample : P0902772-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

0.00min 0.000ng/ml d

response 0

tlc 8/21/09 MP

KR 8/22/09

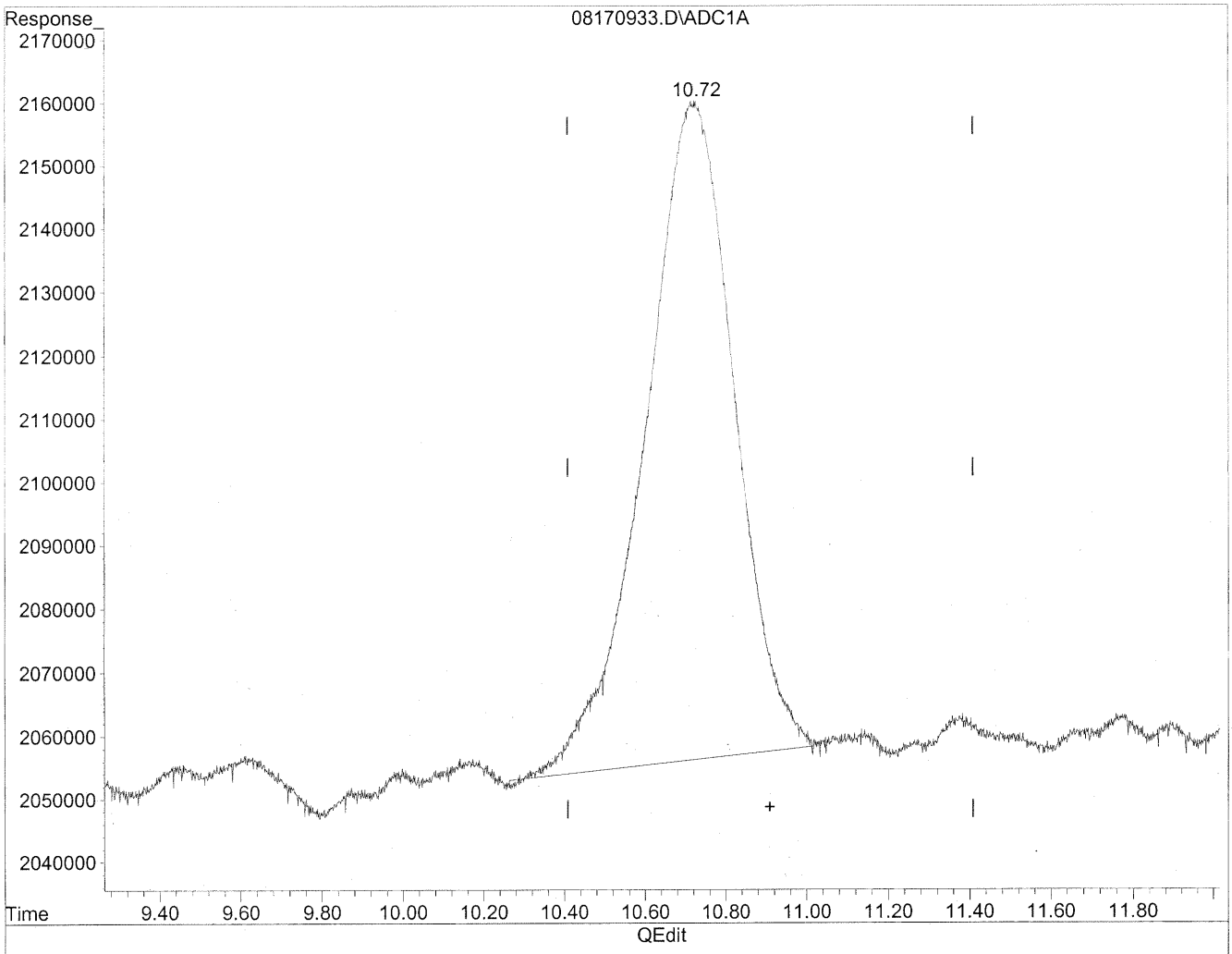
(+) = Expected Retention Time

08170933.D TO110709.M Fri Aug 21 18:14:41 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170933.D Vial: 32
Acq On : 17 Aug 2009 10:51 pm Operator: HC
Sample : P0902772-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.72min 325.151ng/ml

response 15936741

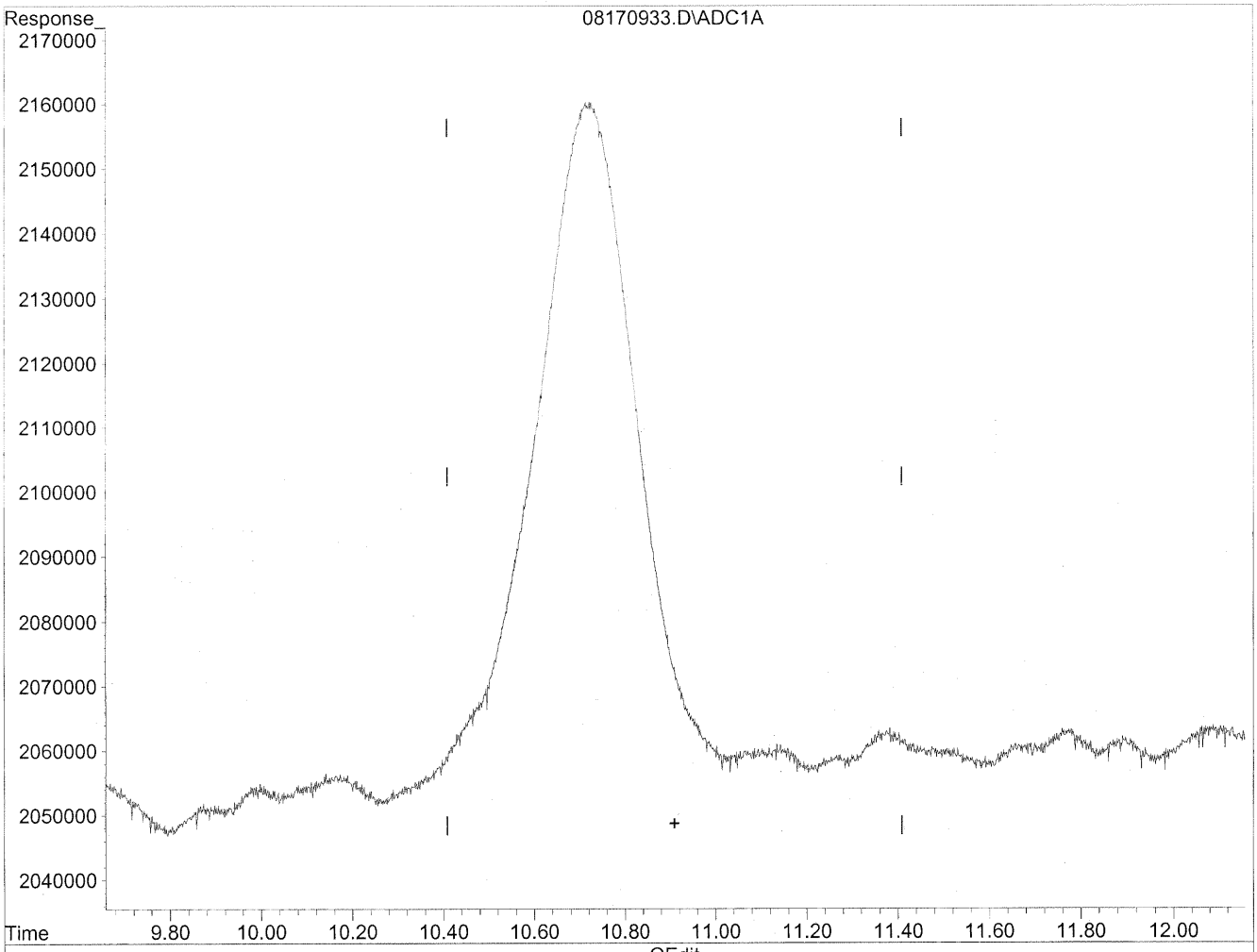
(+) = Expected Retention Time

08170933.D TO110709.M Fri Aug 21 18:14:44 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170933.D Vial: 32
Acq On : 17 Aug 2009 10:51 pm Operator: HC
Sample : P0902772-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

0.00min 0.000ng/ml d

response 0

JL 8/21/09 mp

KP 8/23/09

(+) = Expected Retention Time

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **100308**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P0902772-005

Test Code: EPA Method TO-11A
Instrument ID: Waters LC Module I Plus/UV_Vis 360/LC1
Analyst: Hani Cherazaie
Sampling Media: Silica Gel DNPH Tube
Test Notes: **BC**

Date Collected: 8/10/09
Date Received: 8/12/09
Date Analyzed: 8/17 - 8/18/09
Desorption Volume: 1.0 ml
Volume Sampled: 97.6 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	Formaldehyde	< 100	ND	1.0	ND	0.83	
75-07-0	Acetaldehyde	< 100	ND	1.0	ND	0.57	
123-38-6	Propionaldehyde	< 100	ND	1.0	ND	0.43	
4170-30-3	Crotonaldehyde, Total	< 100	ND	1.0	ND	0.36	
123-72-8	Butyraldehyde	< 100	ND	1.0	ND	0.35	
100-52-7	Benzaldehyde	< 100	ND	1.0	ND	0.24	
590-86-3	Isovaleraldehyde	< 100	ND	1.0	ND	0.29	
110-62-3	Valeraldehyde	< 100	ND	1.0	ND	0.29	
529-20-4	o-Tolualdehyde	< 100	ND	1.0	ND	0.21	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	2.0	ND	0.42	
66-25-1	n-Hexaldehyde	< 100	ND	1.0	ND	0.25	
5779-94-2	**2,5-Dimethylbenzaldehyde**	**730**	**7.5**	1.0	**1.4**	0.19	

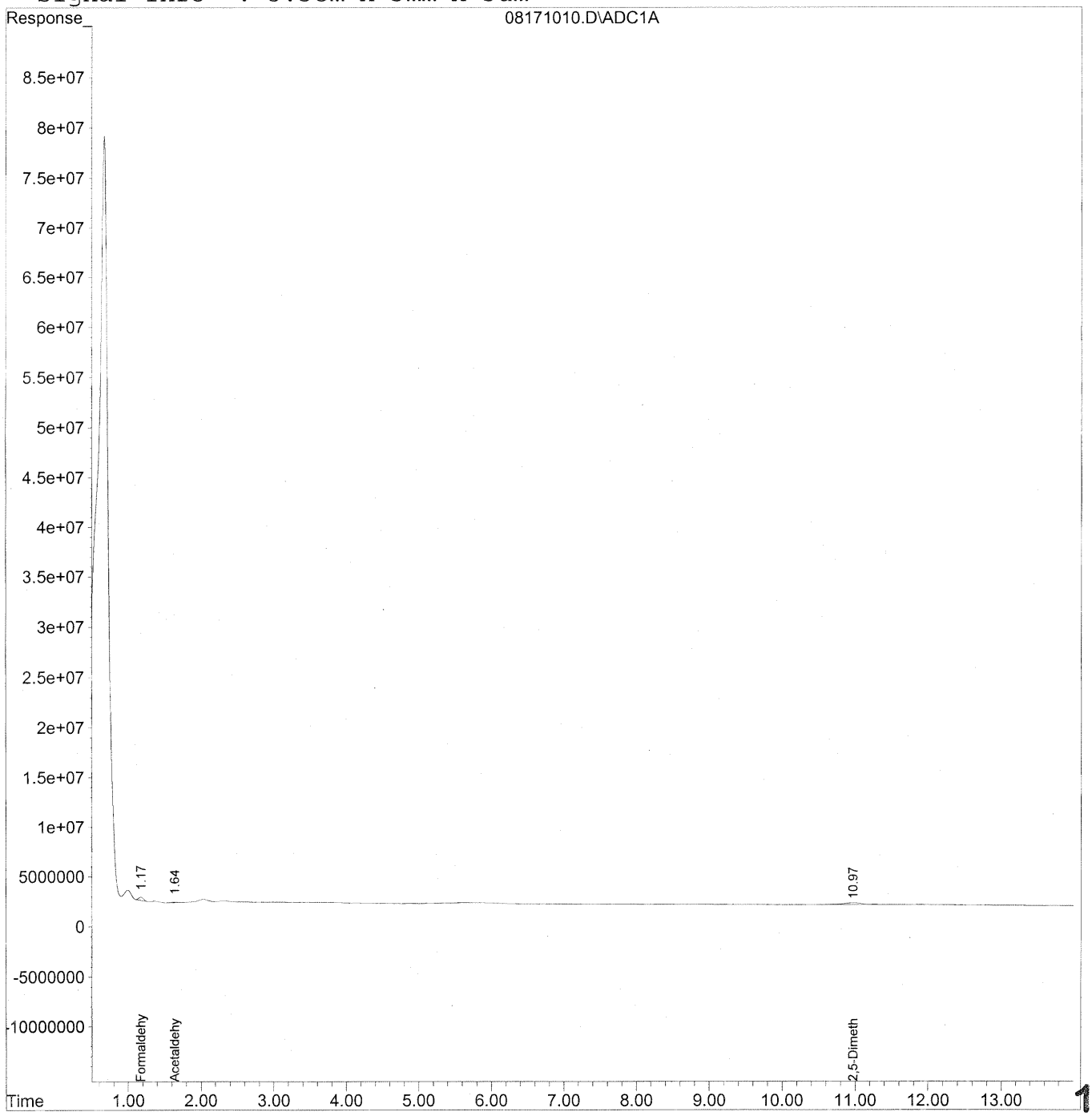
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.

Verified By: _____ Date: 8/25/09

Data File : J:\LC01\DATA\TO11\2009_08\17\08171010.D Vial: 23
Acq On : 18 Aug 2009 6:09 pm Operator: HC
Sample : P0902772-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:22 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



120

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171010.D Vial: 23
Acq On : 18 Aug 2009 6:09 pm Operator: HC
Sample : P0902772-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:22 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

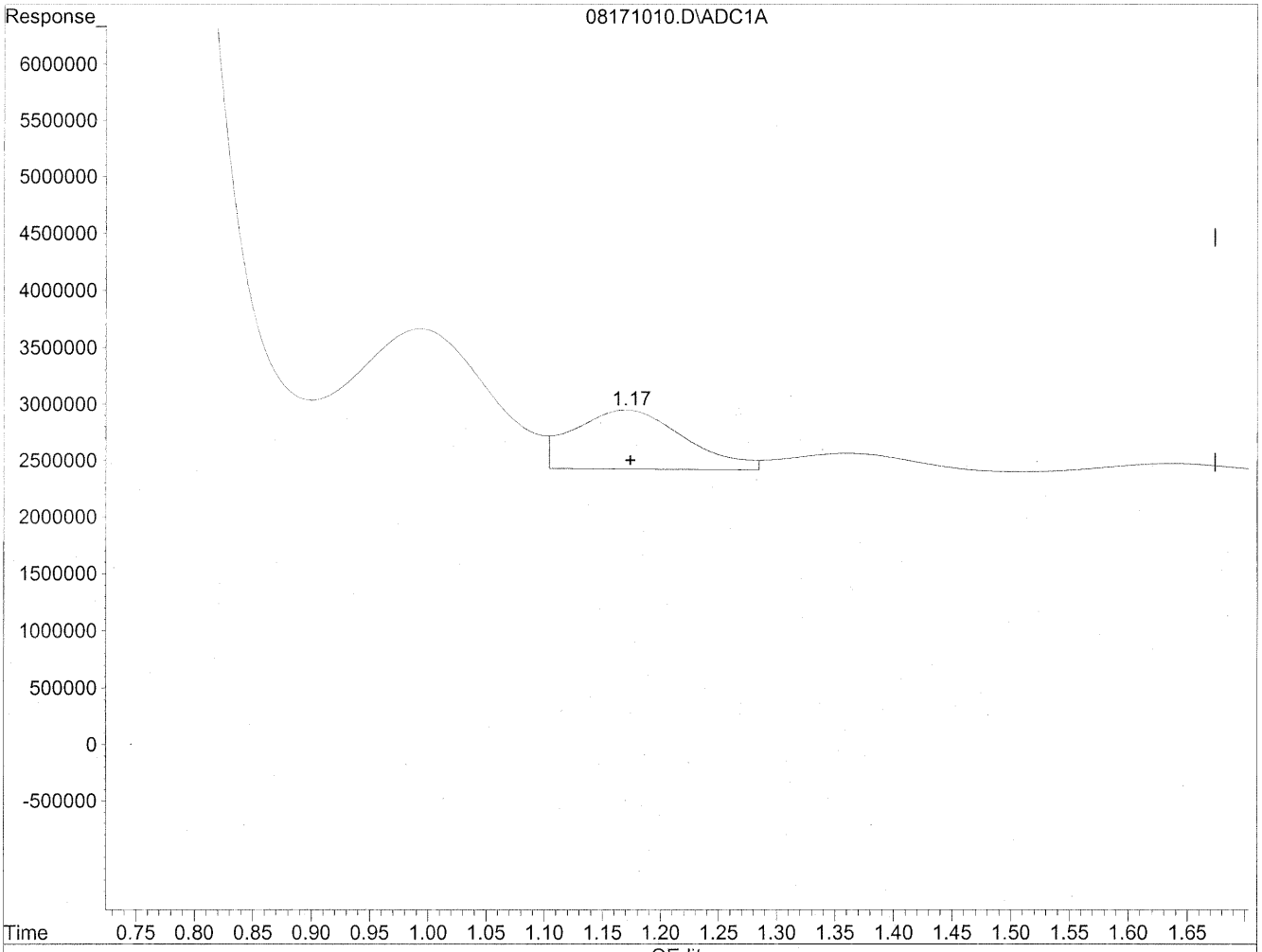
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	14846307	80.870	ng/mlm
2)	Acetaldehyde	1.64	4704196	33.548	ng/ml
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/mld
12)	2,5-Dimethylbenzaldehyde	10.97	35876312	731.970	ng/mlm

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171010.D Vial: 23
Acq On : 18 Aug 2009 6:09 pm Operator: HC
Sample : P0902772-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.17min 190.218ng/ml

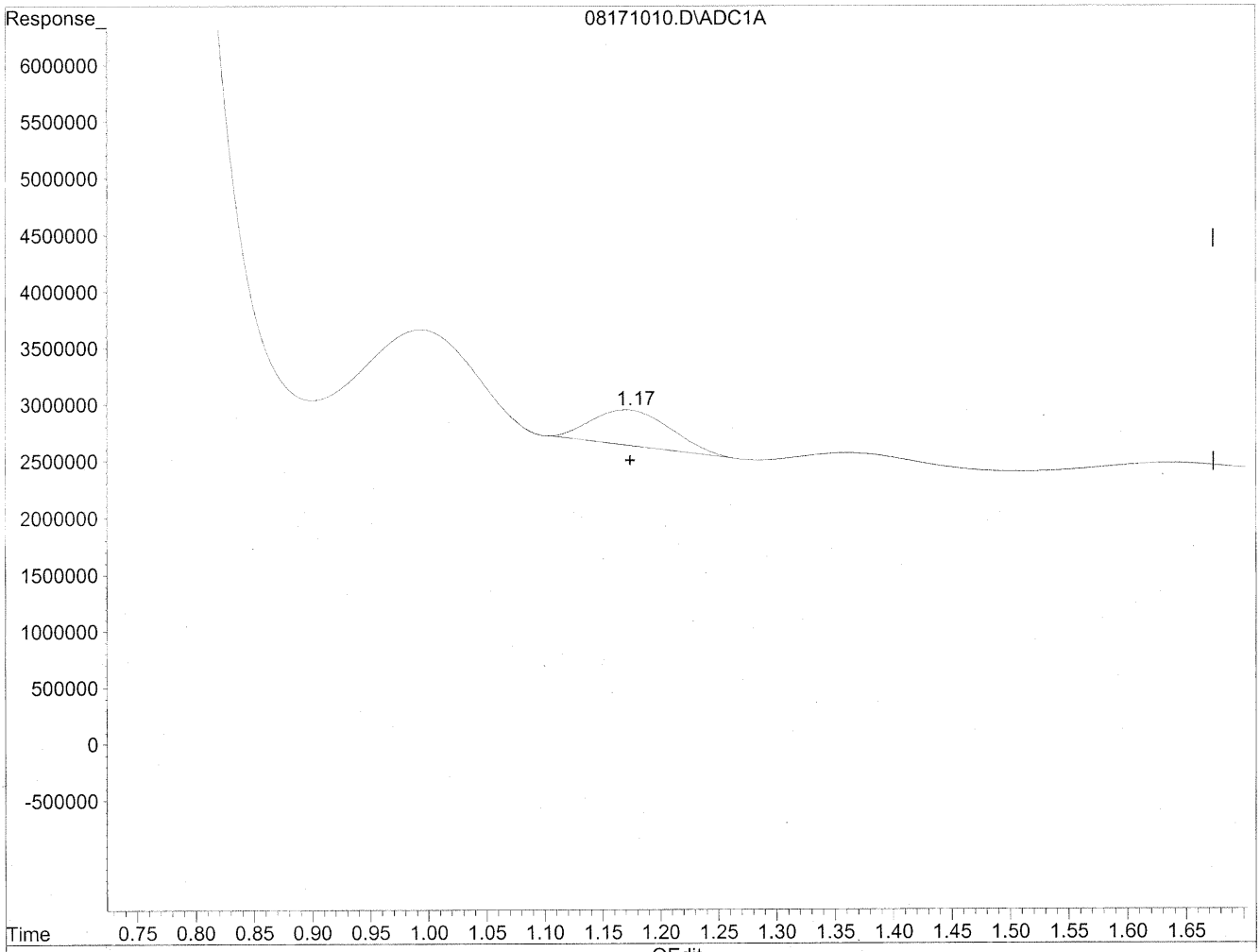
response 34920549

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171010.D Vial: 23
Acq On : 18 Aug 2009 6:09 pm Operator: HC
Sample : P0902772-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 80.870ng/ml m

response 14846307

HC 8/22/09

BC

KP 8/23/09

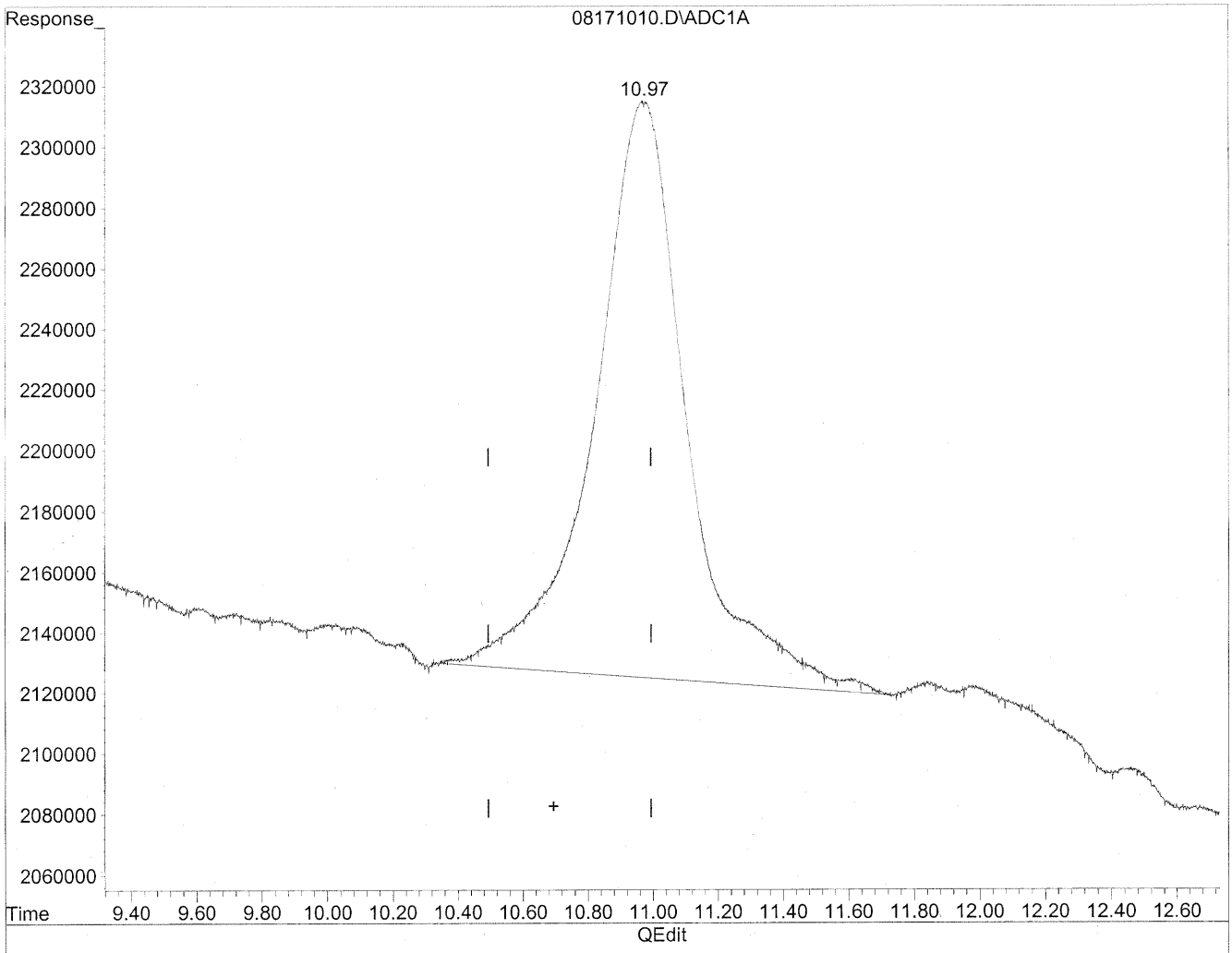
(+) = Expected Retention Time

08171010.D TO110709.M Sat Aug 22 14:21:46 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171010.D Vial: 23
Acq On : 18 Aug 2009 6:09 pm Operator: HC
Sample : P0902772-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.97min 557.565ng/ml

response 37548519

124

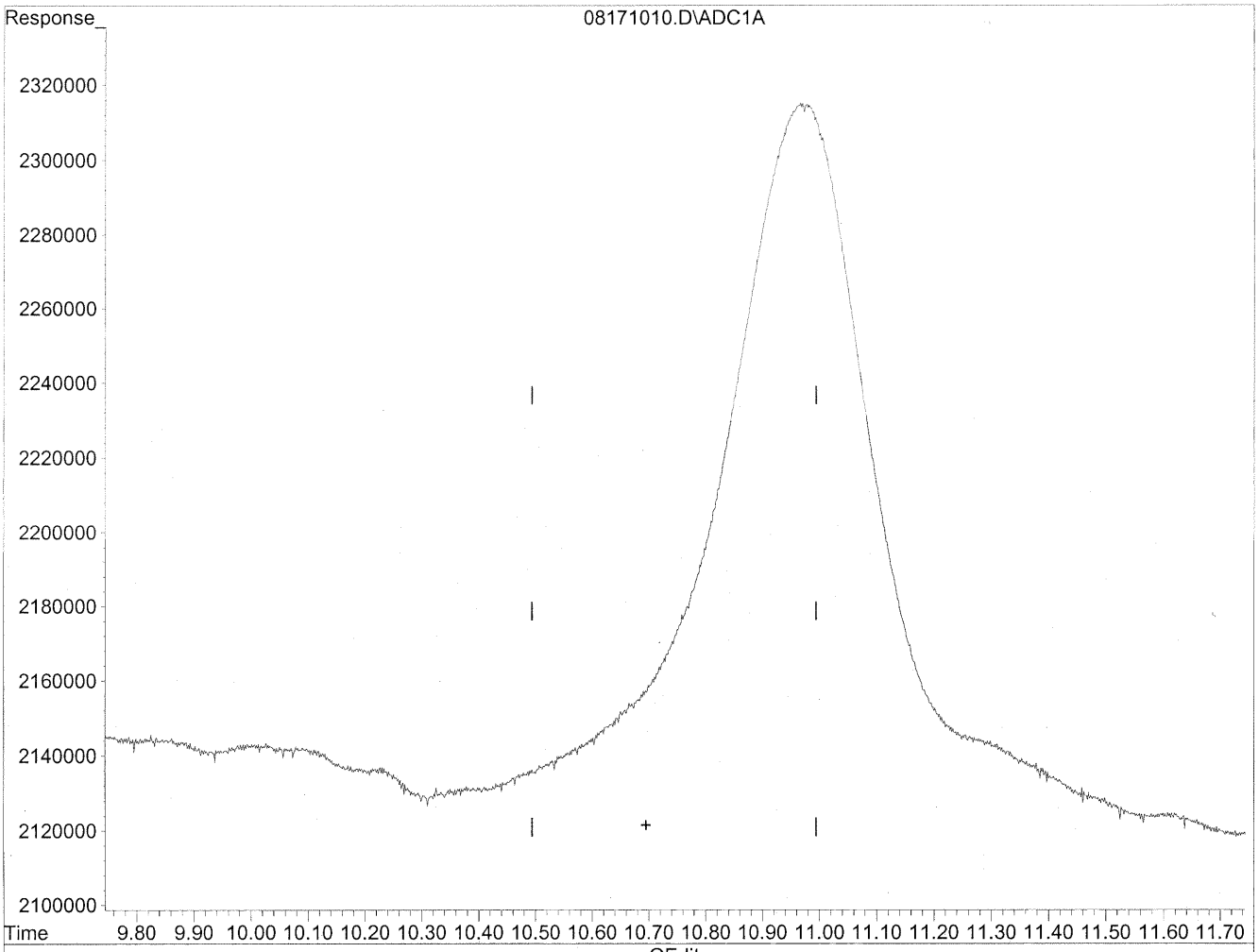
(+) = Expected Retention Time

08171010.D TO110709.M Sat Aug 22 14:21:59 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171010.D Vial: 23
Acq On : 18 Aug 2009 6:09 pm Operator: HC
Sample : P0902772-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09 WP 8/23/09

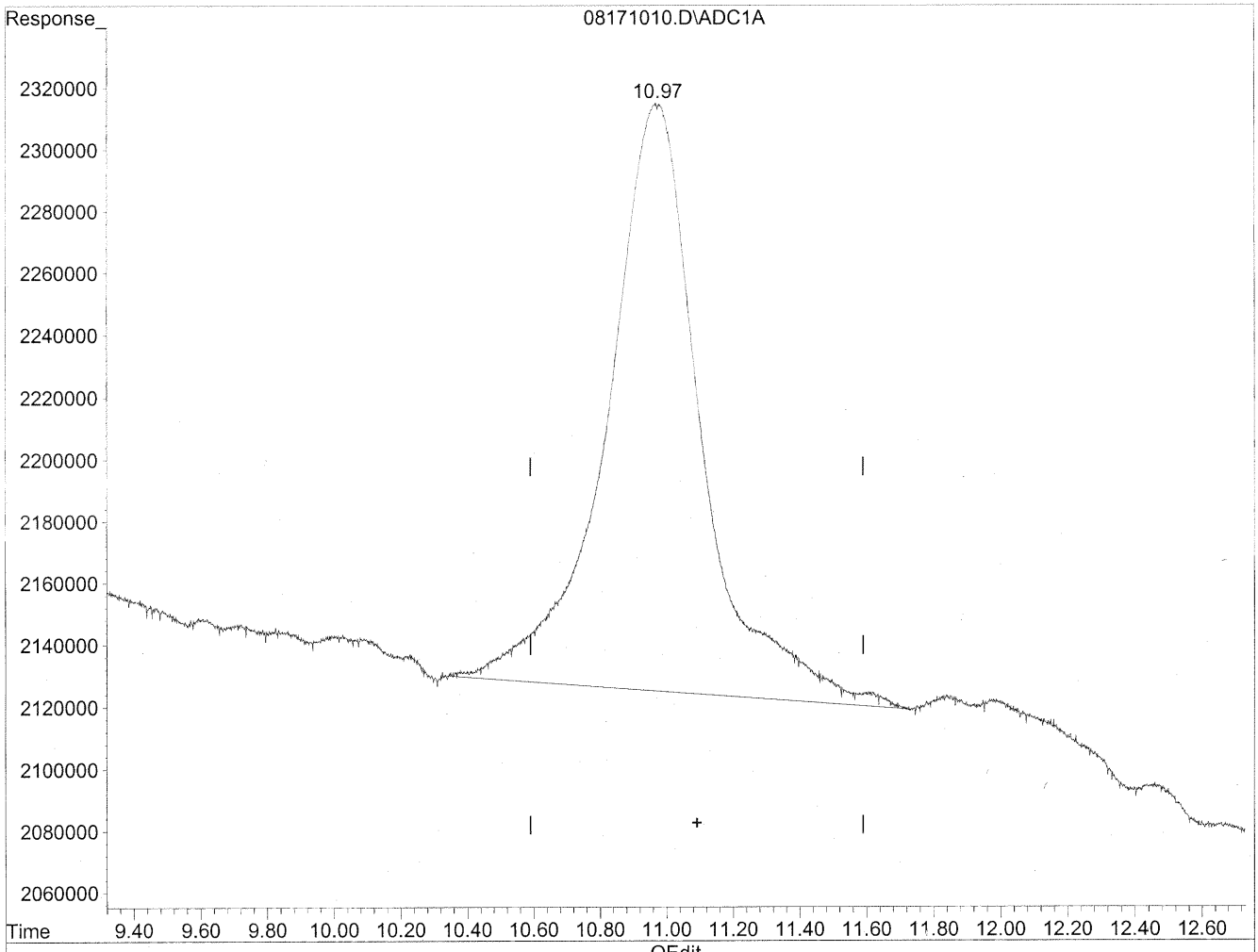
(+) = Expected Retention Time

08171010.D TO110709.M Sat Aug 22 14:22:02 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171010.D Vial: 23
Acq On : 18 Aug 2009 6:09 pm Operator: HC
Sample : P0902772-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(12) 2,5-Dimethylbenzaldehyde

10.97min 766.087ng/ml

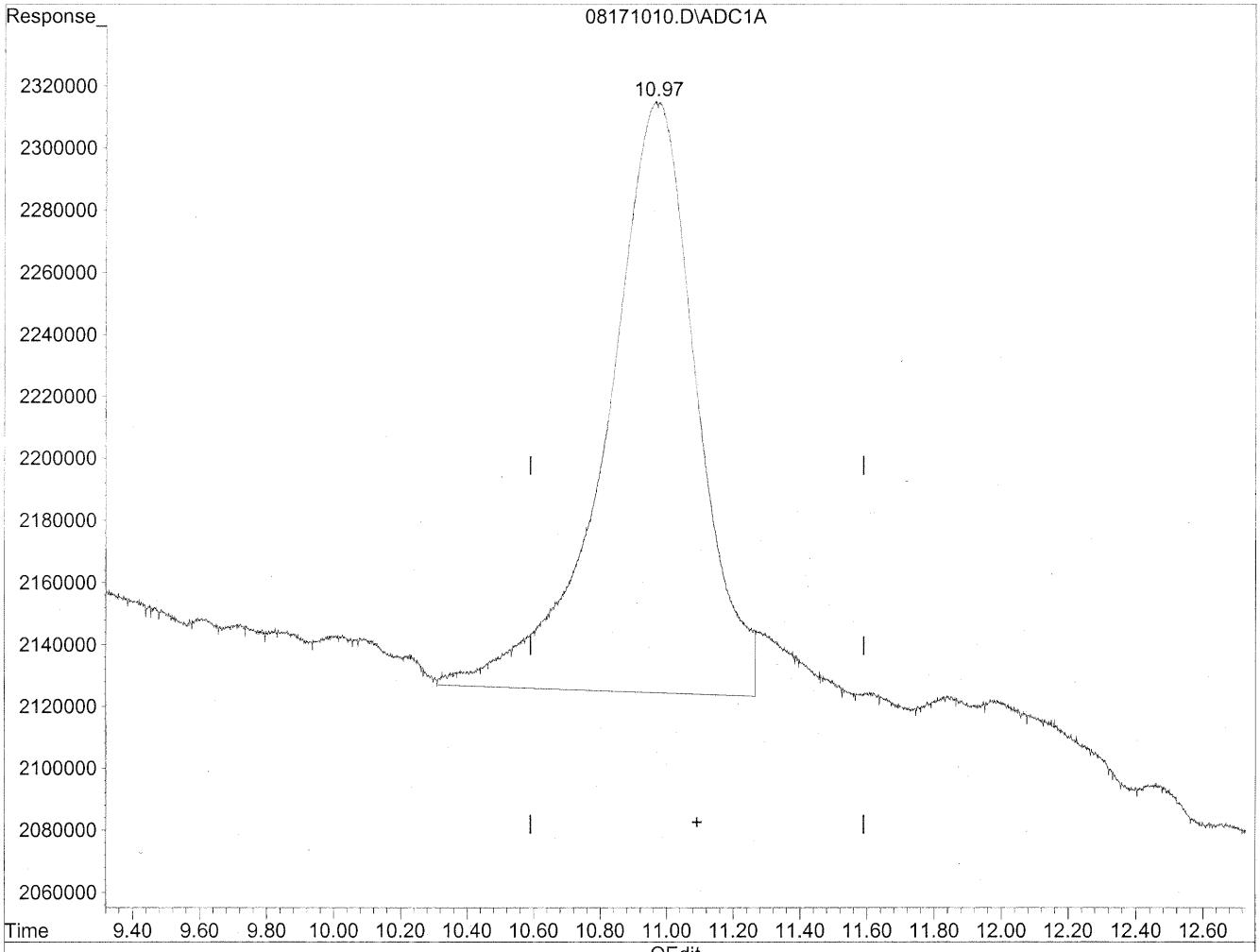
response 37548519

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171010.D Vial: 23
Acq On : 18 Aug 2009 6:09 pm Operator: HC
Sample : P0902772-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.97min 731.970ng/ml m

response 35876312

HC 8/22/09

IC KR 8/23/09

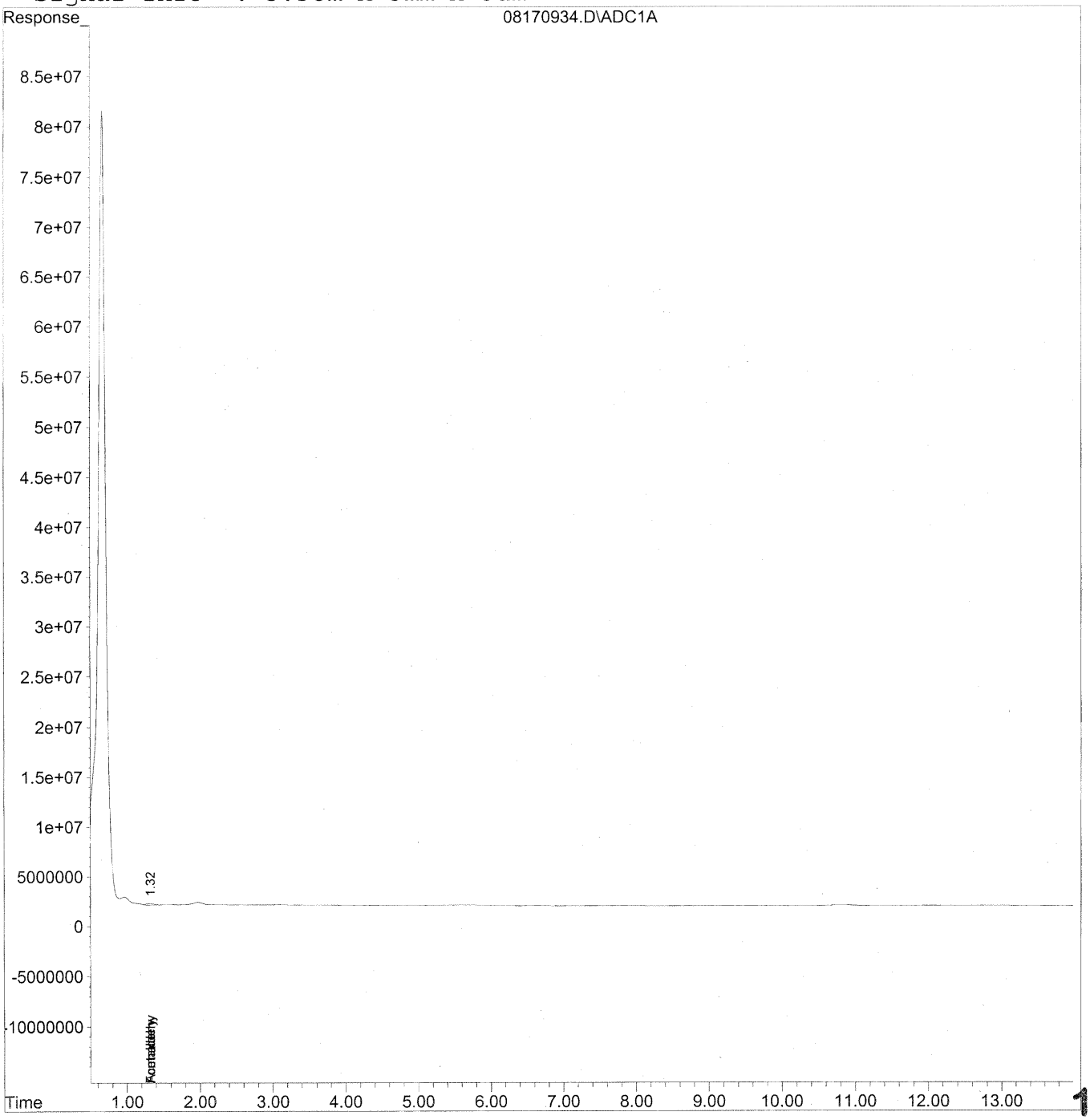
(+) = Expected Retention Time

08171010.D TO110709.M Sat Aug 22 14:22:13 2009

Data File : J:\LC01\DATA\TO11\2009_08\17\08170934.D Vial: 33
Acq On : 17 Aug 2009 11:06 pm Operator: HC
Sample : P0902772-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:15 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170934.D Vial: 33
Acq On : 17 Aug 2009 11:06 pm Operator: HC
Sample : P0902772-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:15 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

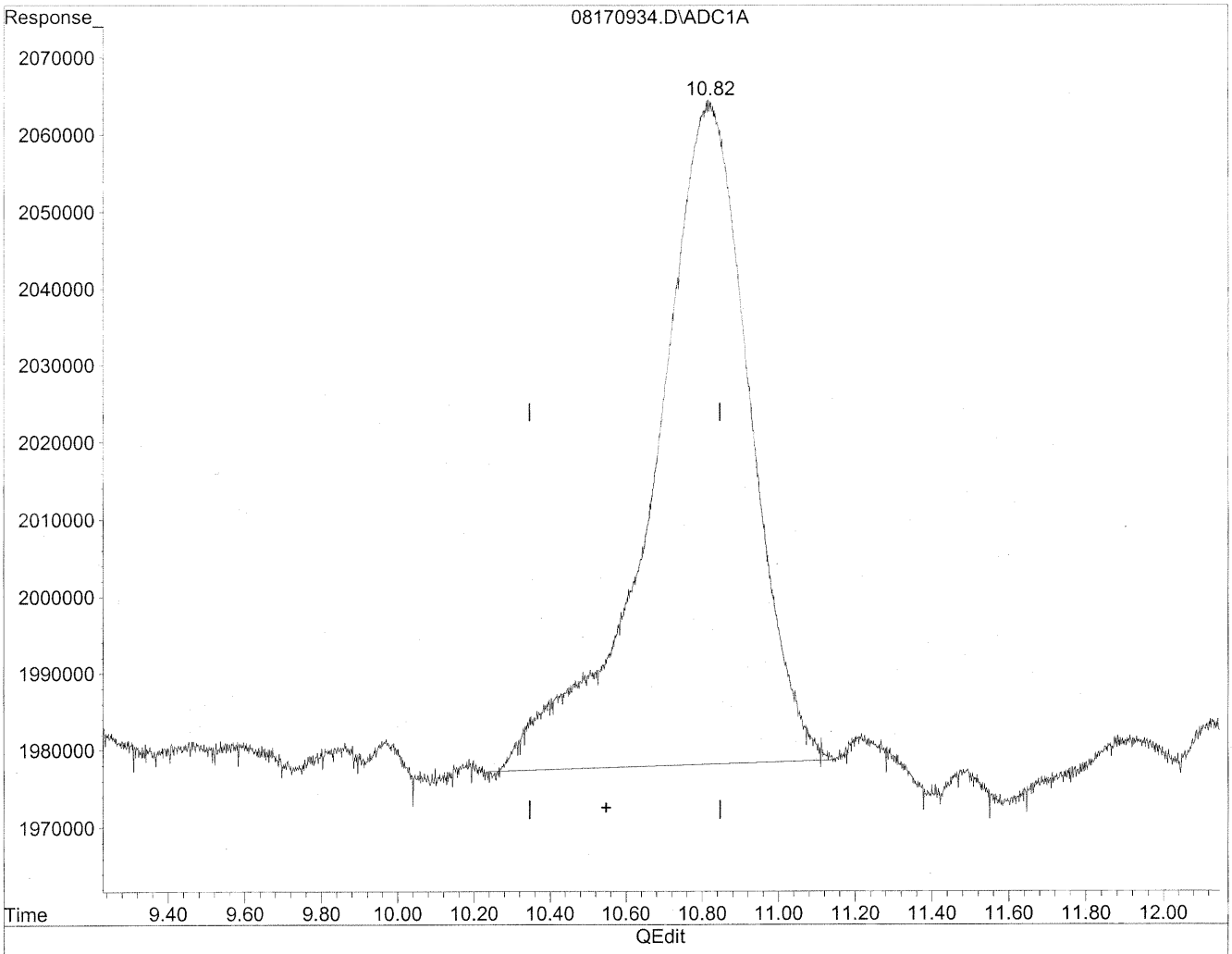
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.32	11043166	60.154	ng/ml
2)	Acetaldehyde	1.32f	11043166	78.754	ng/ml
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/mld
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/mld

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170934.D Vial: 33
Acq On : 17 Aug 2009 11:06 pm Operator: HC
Sample : P0902772-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:15 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.82min 225.594ng/ml

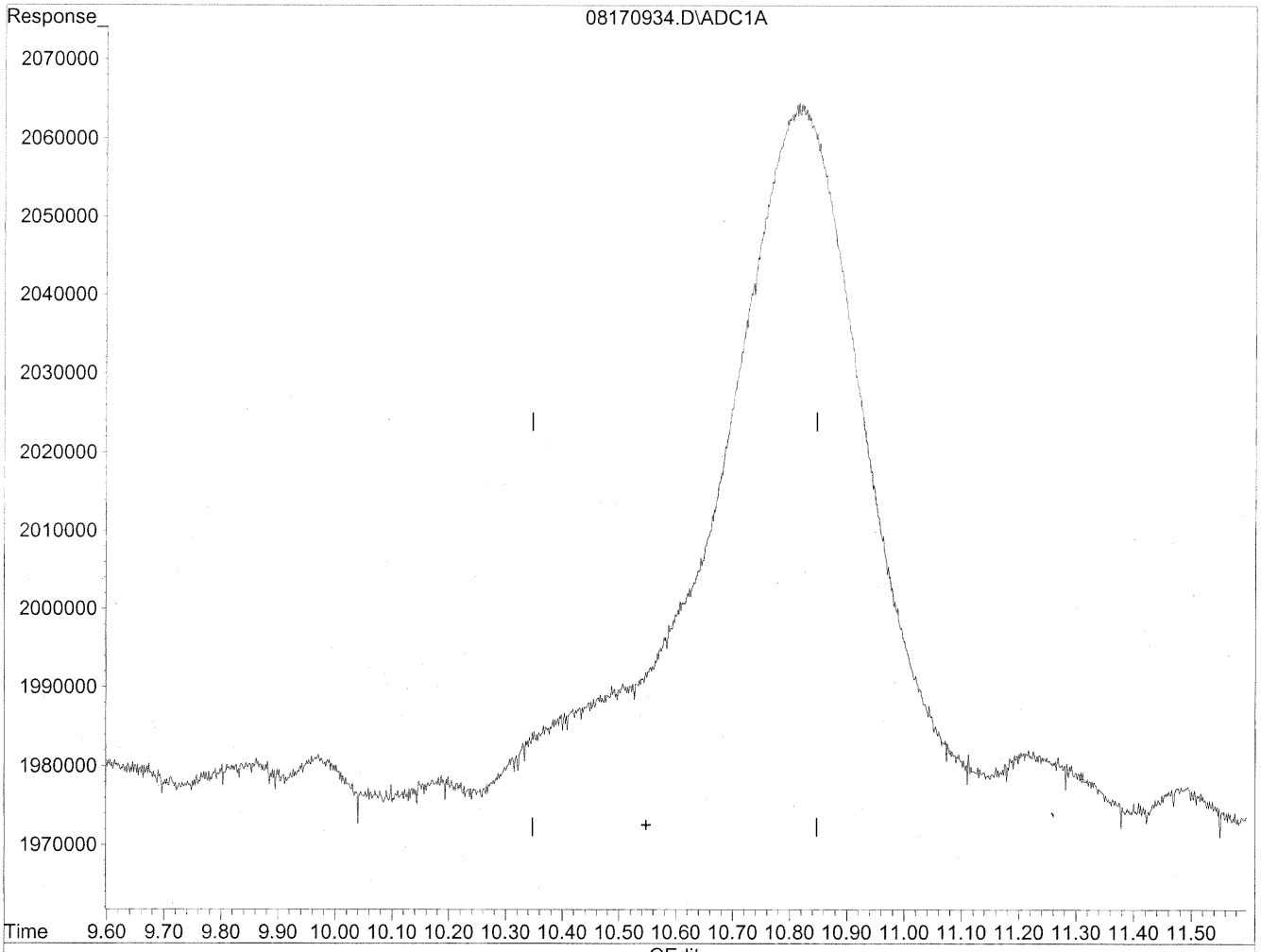
response 15192320

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170934.D Vial: 33
Acq On : 17 Aug 2009 11:06 pm Operator: HC
Sample : P0902772-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:15 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/24/09

KP 8/23/09

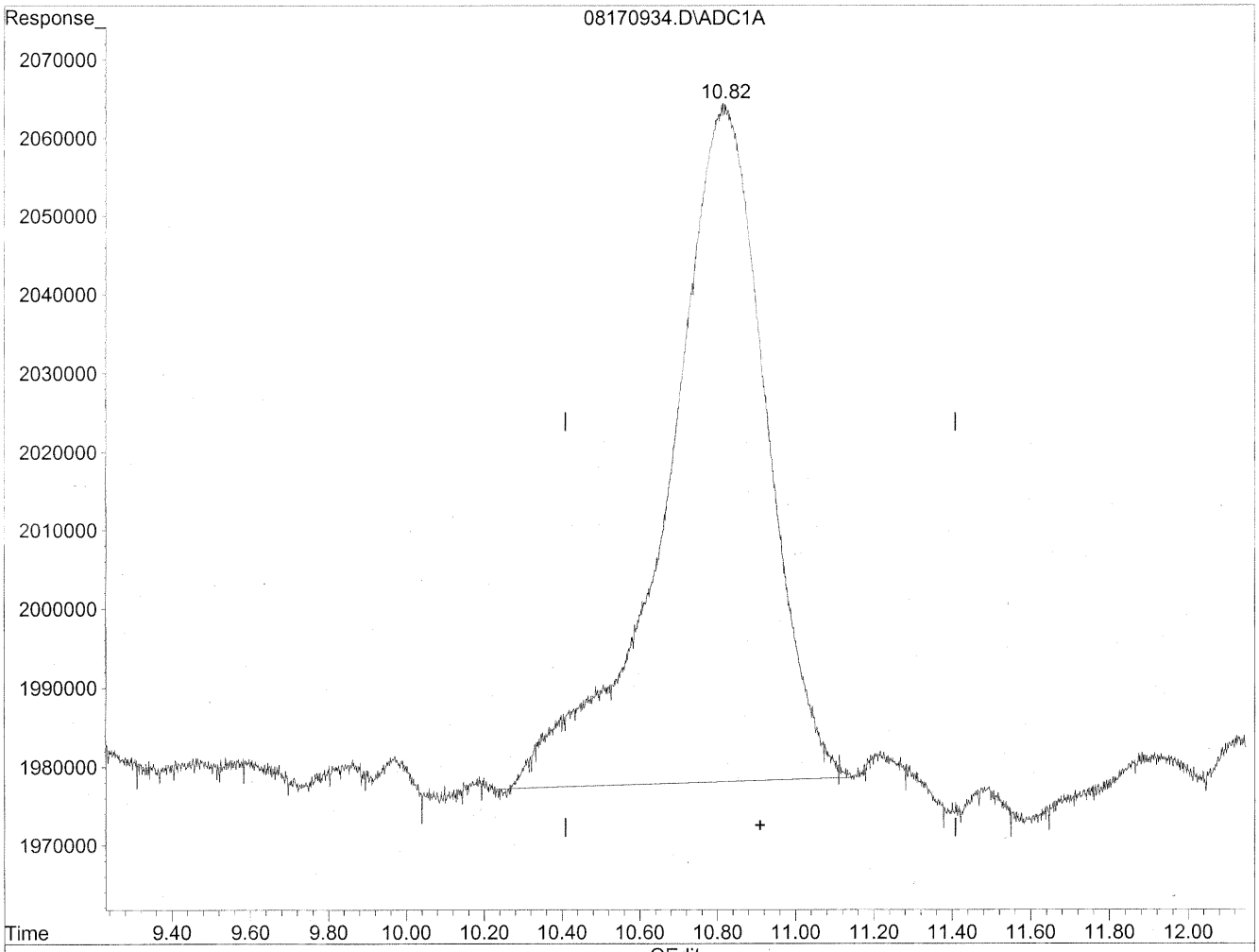
(+) = Expected Retention Time

08170934.D TO110709.M Fri Aug 21 18:15:29 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170934.D Vial: 33
Acq On : 17 Aug 2009 11:06 pm Operator: HC
Sample : P0902772-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:15 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.82min 309.963ng/ml

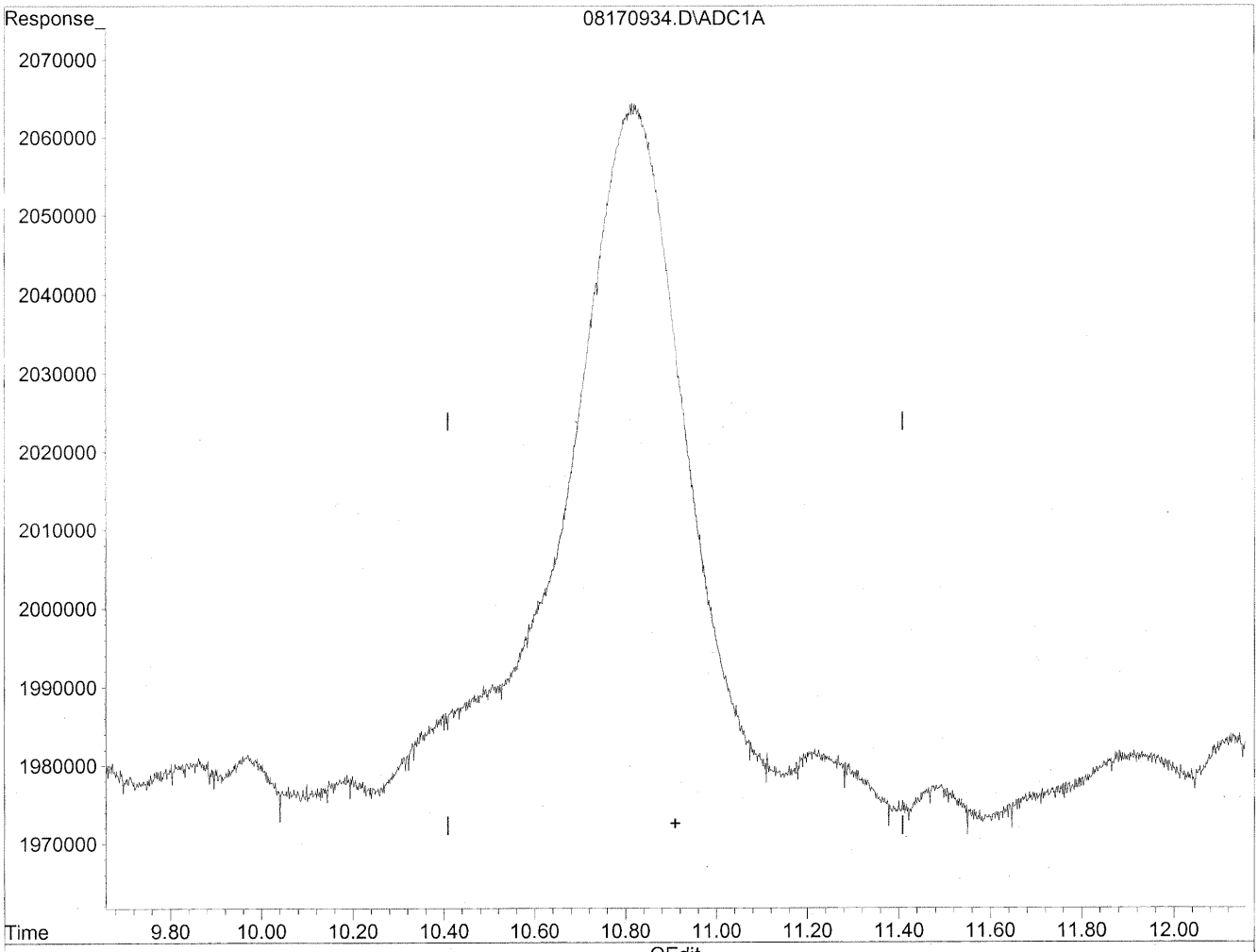
response 15192320

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170934.D Vial: 33
Acq On : 17 Aug 2009 11:06 pm Operator: HC
Sample : P0902772-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:15 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(12) 2,5-Dimethylbenzaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/21/09
MP
KP 8/23/09

133

(+) = Expected Retention Time

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **100316**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P0902772-006

Test Code: EPA Method TO-11A
Instrument ID: Waters LC Module I Plus/UV_Vis 360/LC1
Analyst: Hani Cherazaie
Sampling Media: Silica Gel DNPH Tube
Test Notes: **BC**

Date Collected: 8/10/09
Date Received: 8/12/09
Date Analyzed: 8/17 - 8/18/09
Desorption Volume: 1.0 ml
Volume Sampled: 107.1 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**4,000**	**37**	0.93	**30**	0.76	
75-07-0	**Acetaldehyde**	**2,600**	**24**	0.93	**13**	0.52	**BT**
123-38-6	**Propionaldehyde**	**230**	**2.2**	0.93	**0.92**	0.39	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.93	ND	0.33	
123-72-8	**Butyraldehyde**	**430**	**4.0**	0.93	**1.4**	0.32	**M**
100-52-7	**Benzaldehyde**	**410**	**3.8**	0.93	**0.88**	0.22	
590-86-3	Isovaleraldehyde	< 100	ND	0.93	ND	0.27	
110-62-3	**Valeraldehyde**	**380**	**3.6**	0.93	**1.0**	0.27	
529-20-4	o-Tolualdehyde	< 100	ND	0.93	ND	0.19	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	1.9	ND	0.38	
66-25-1	**n-Hexaldehyde**	**620**	**5.8**	0.93	**1.4**	0.23	
5779-94-2	**2,5-Dimethylbenzaldehyde**	**550**	**5.1**	0.93	**0.93**	0.17	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
BT = Results indicated possible breakthrough; back section > 10% front section.
M = Matrix interference; results may be biased high.

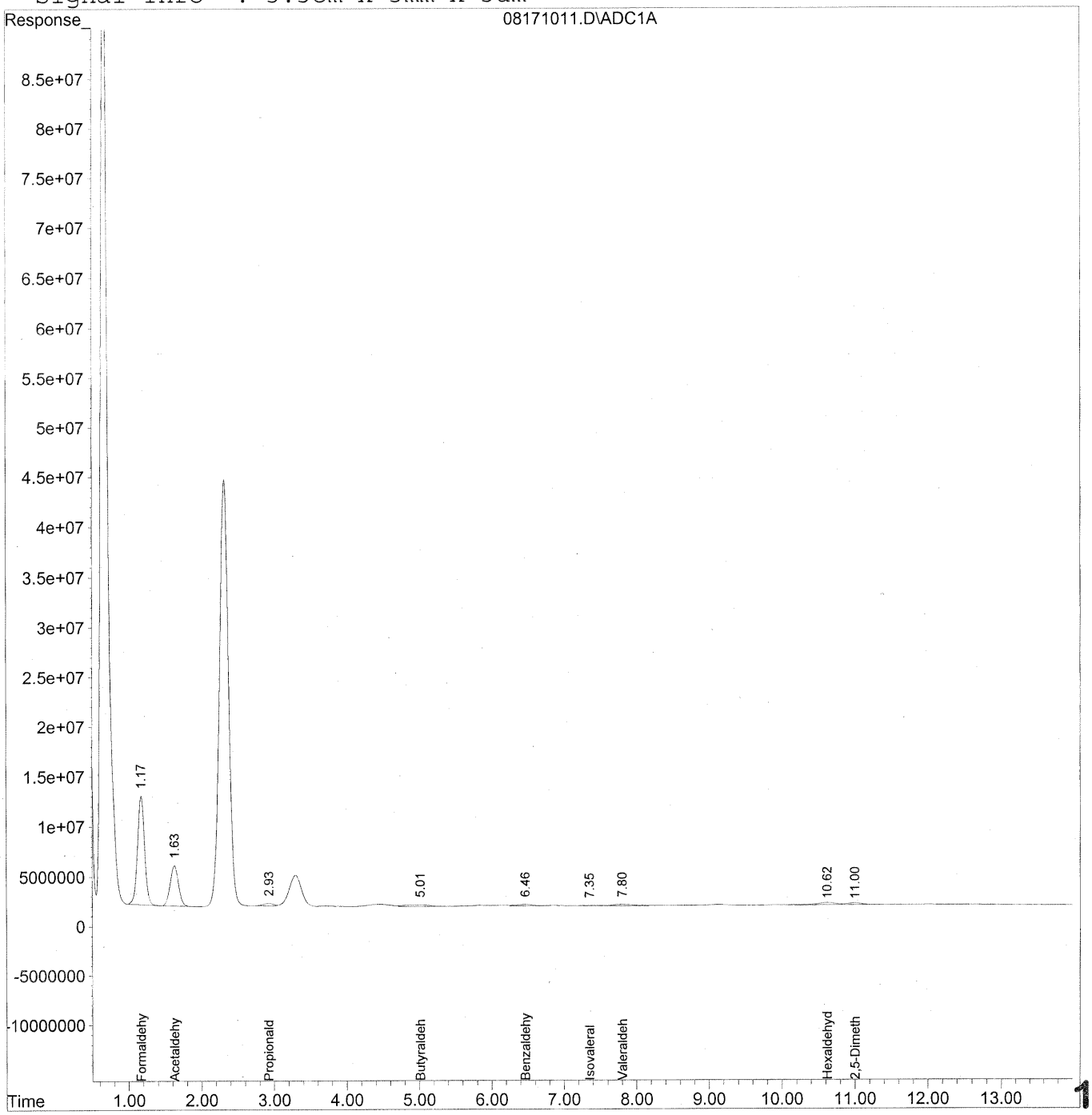
Verified By: _____ Date: 8/25/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:29 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:29 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

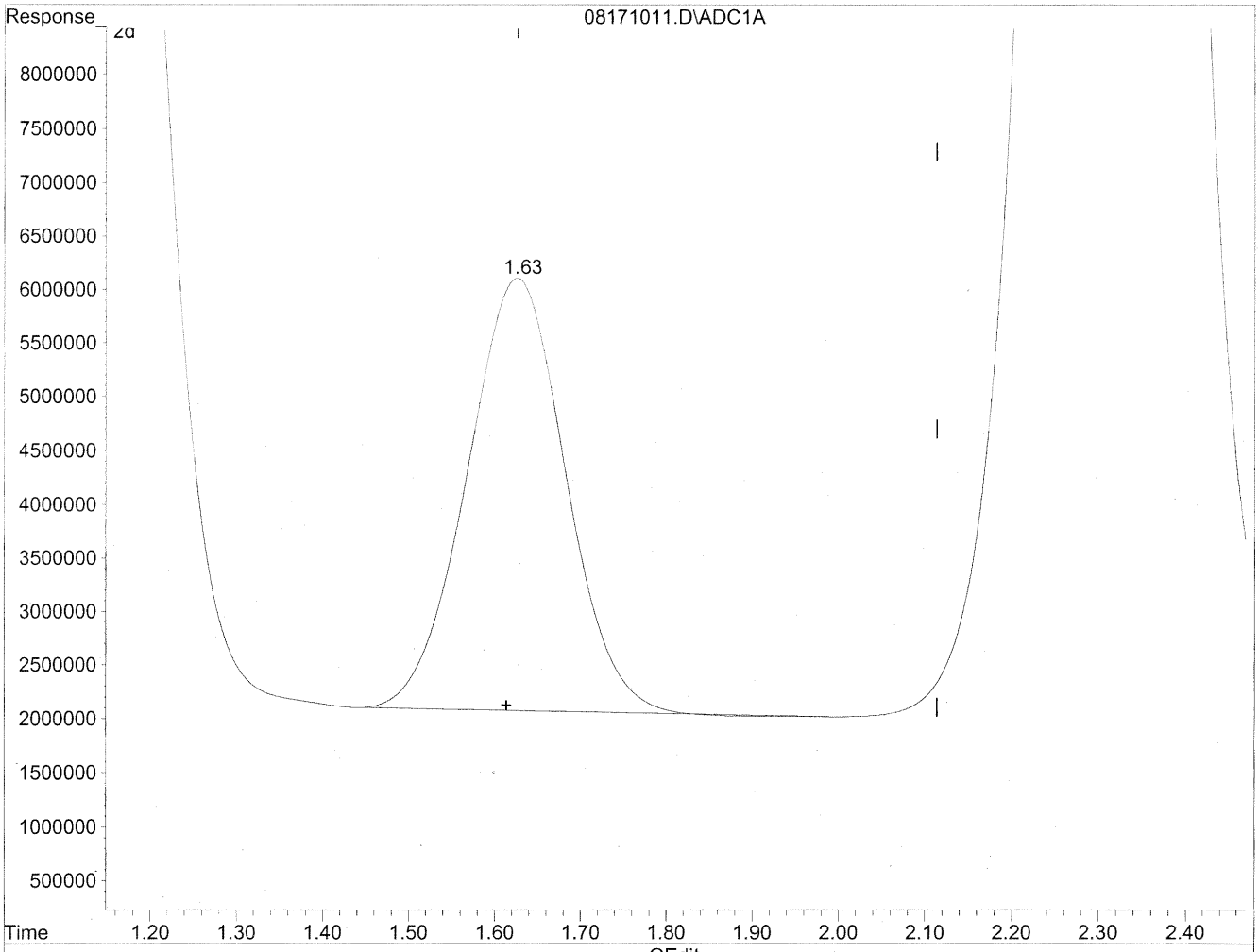
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	732784823	3991.607	ng/ml
2)	Acetaldehyde	1.63	323260097	2305.320	ng/mlm
3)	Propionaldehyde	2.93	24874069	233.132	ng/mlm
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	5.01	37668543	426.423	ng/mlm
6)	Benzaldehyde	6.46	26962027	409.326	ng/mlm
7)	Isovaleraldehyde	7.35	6897046	88.140	ng/mlm
8)	Valeraldehyde	7.80	27999295	380.917	ng/mlm
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/mld
11)	Hexaldehyde	10.62	42025868	624.050	ng/mlm
12)	2,5-Dimethylbenzaldehyde	11.00	26824624	547.292	ng/mlm

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.63min 2291.474ng/ml

response 321318482

137

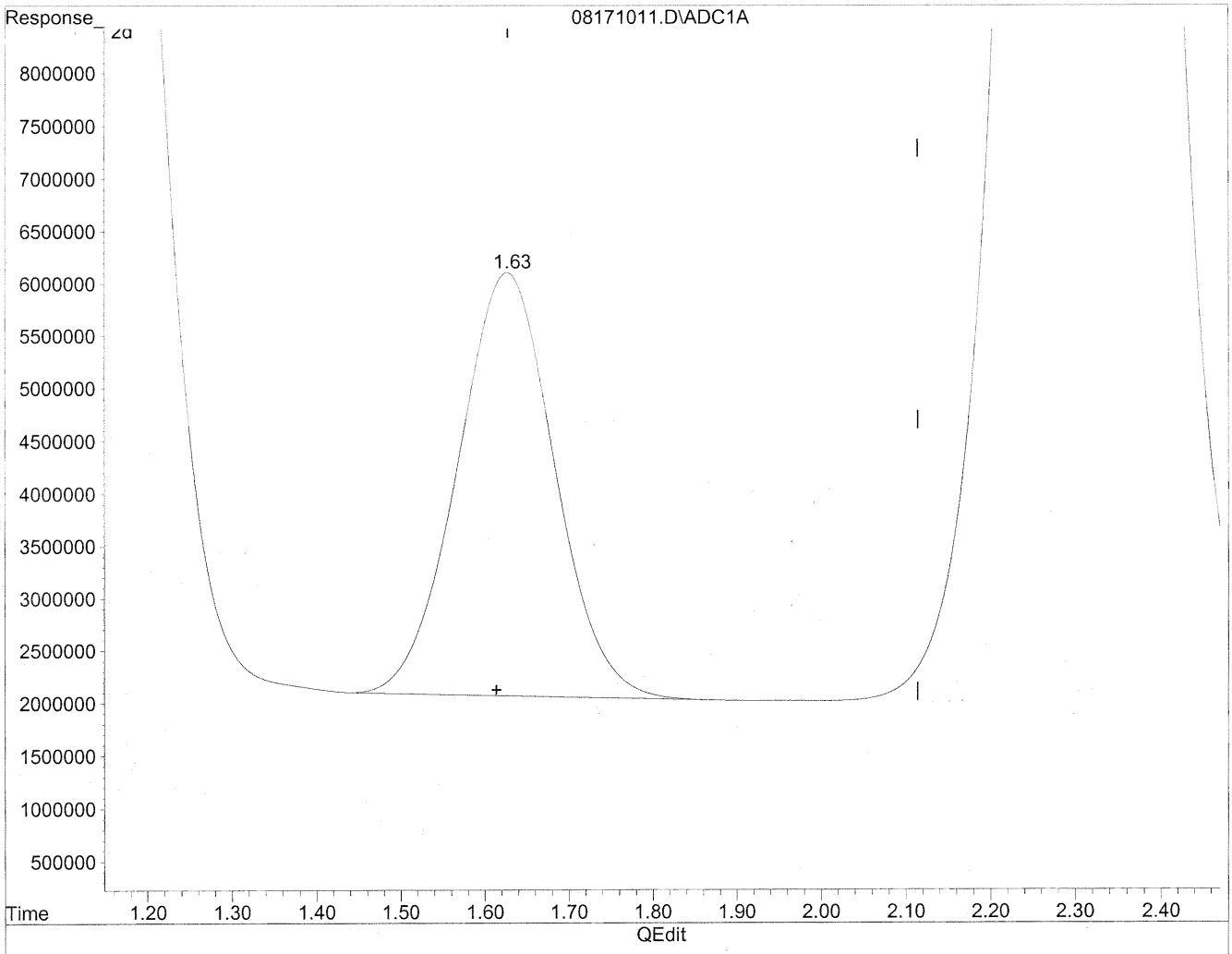
(+) = Expected Retention Time

08171011.D TO110709.M Sat Aug 22 14:27:07 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.63min 2305.320ng/ml m

response 323260097

HC 8/22/09 1C

KP 8/23/09

138

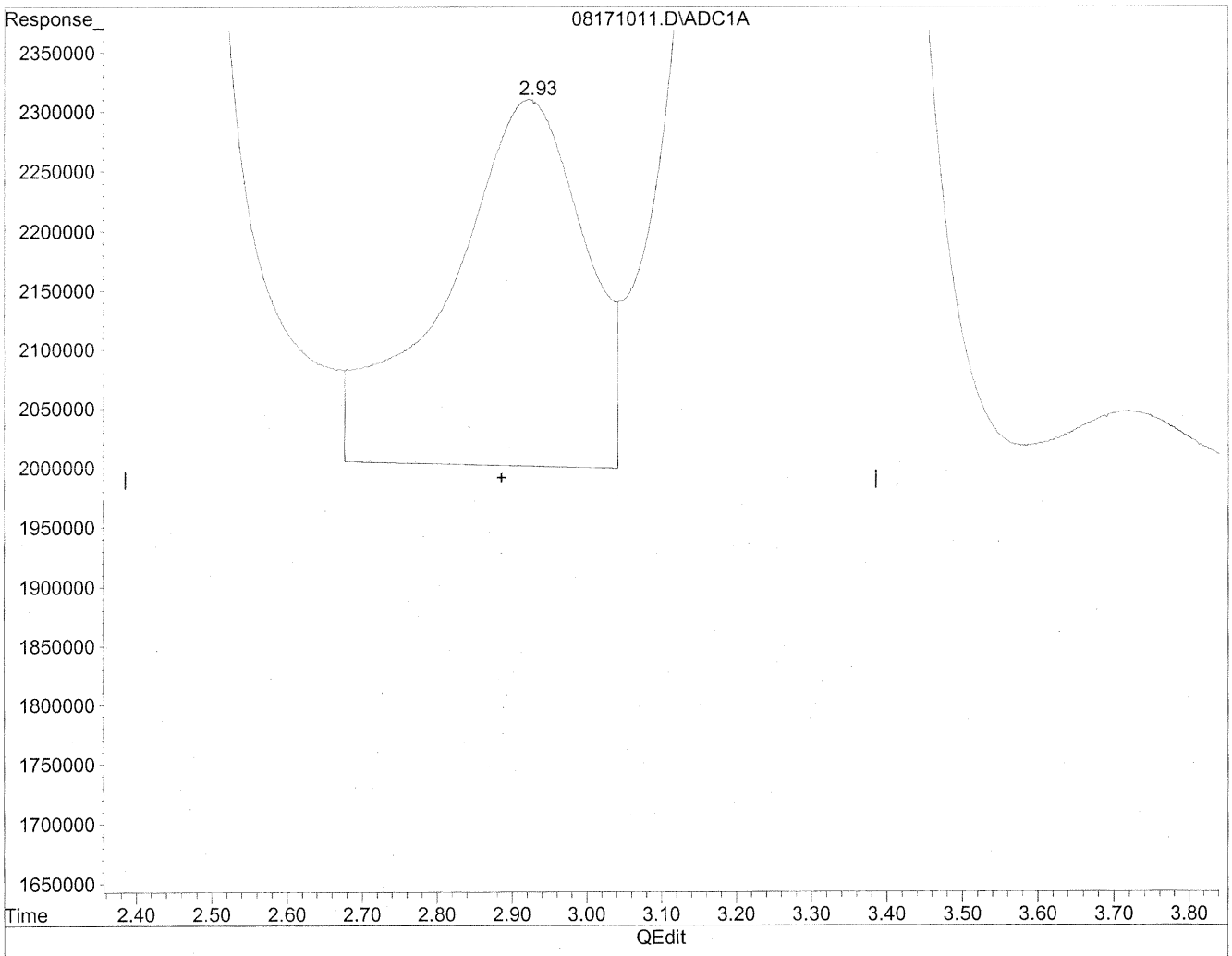
(+) = Expected Retention Time

08171011.D TO110709.M Sat Aug 22 14:27:14 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.92min 368.392ng/ml

response 39305662

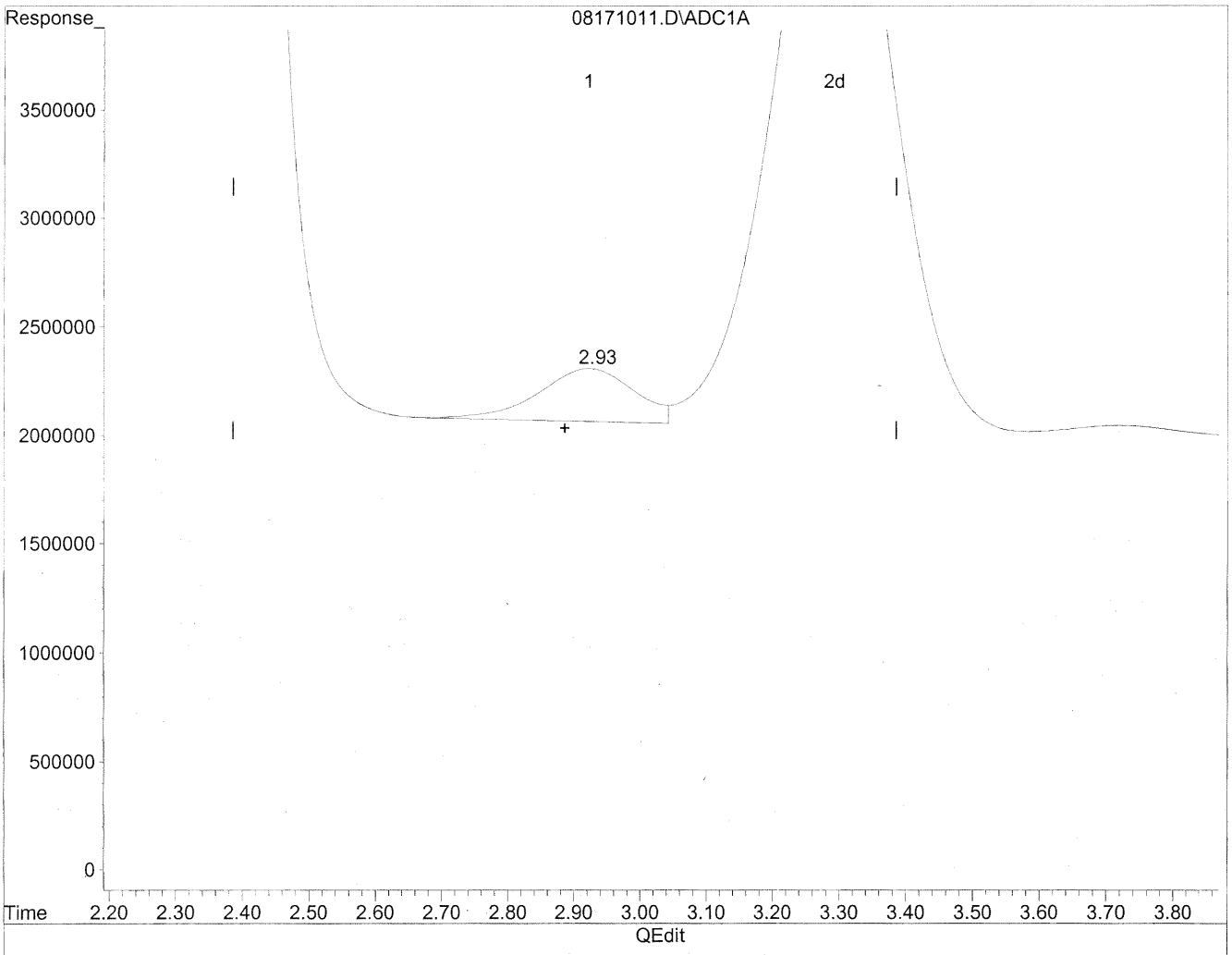
139

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.93min 233.132ng/ml m

response 24874069

HC 8/22/09

BC

KP 8/22/09

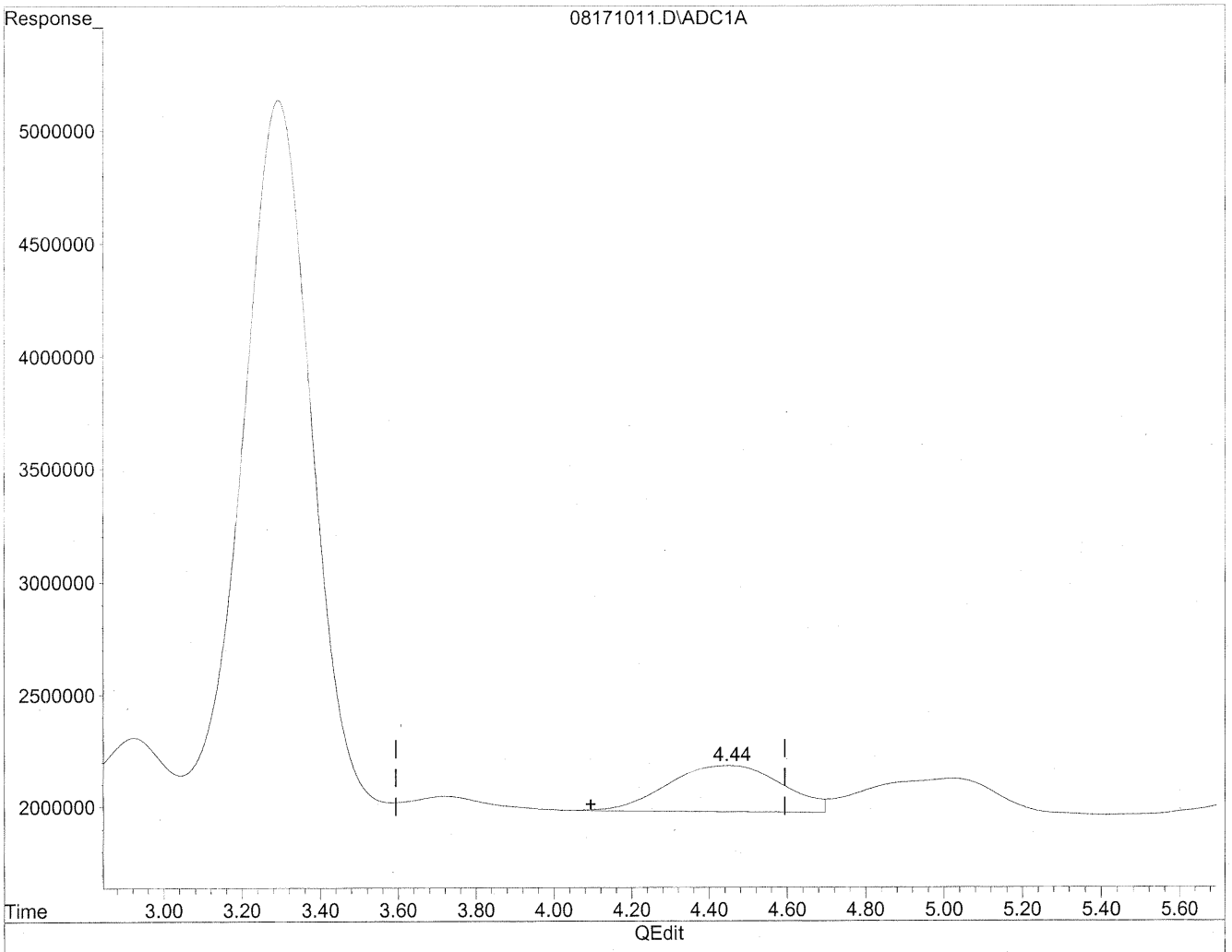
140

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.45min 429.622ng/ml

response 41851749

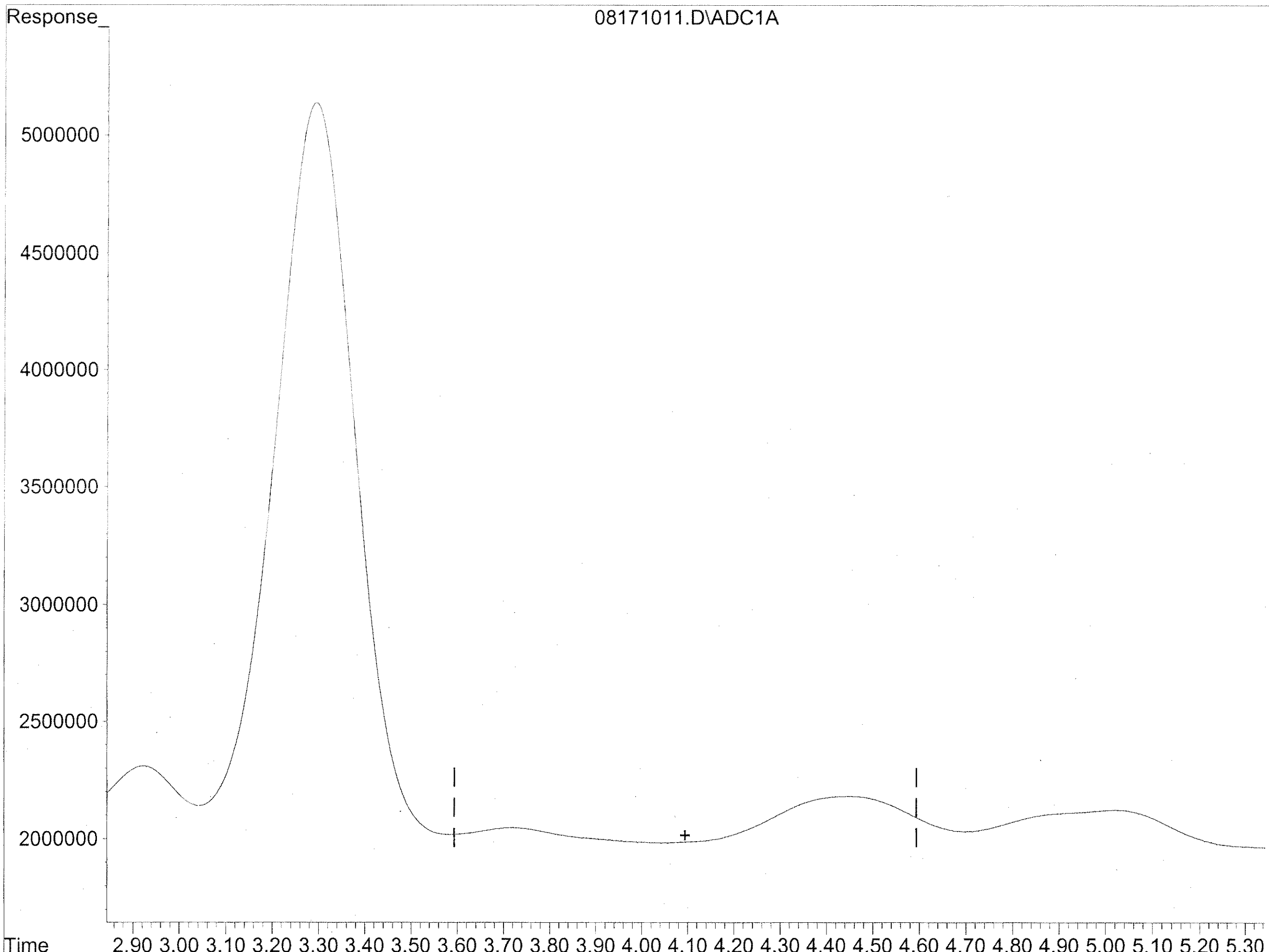
141

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D   Vial: 24
Acq On   : 18 Aug 2009   6:24 pm   Operator: HC
Sample   : P0902772-006 front 1.0ml   Inst   : LC 01
Misc   :   Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method   : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title   : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update   : Wed Aug 19 10:45:48 2009
Response via   : Multiple Level Calibration



QEdit

(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09
MP
KR 8/23/09

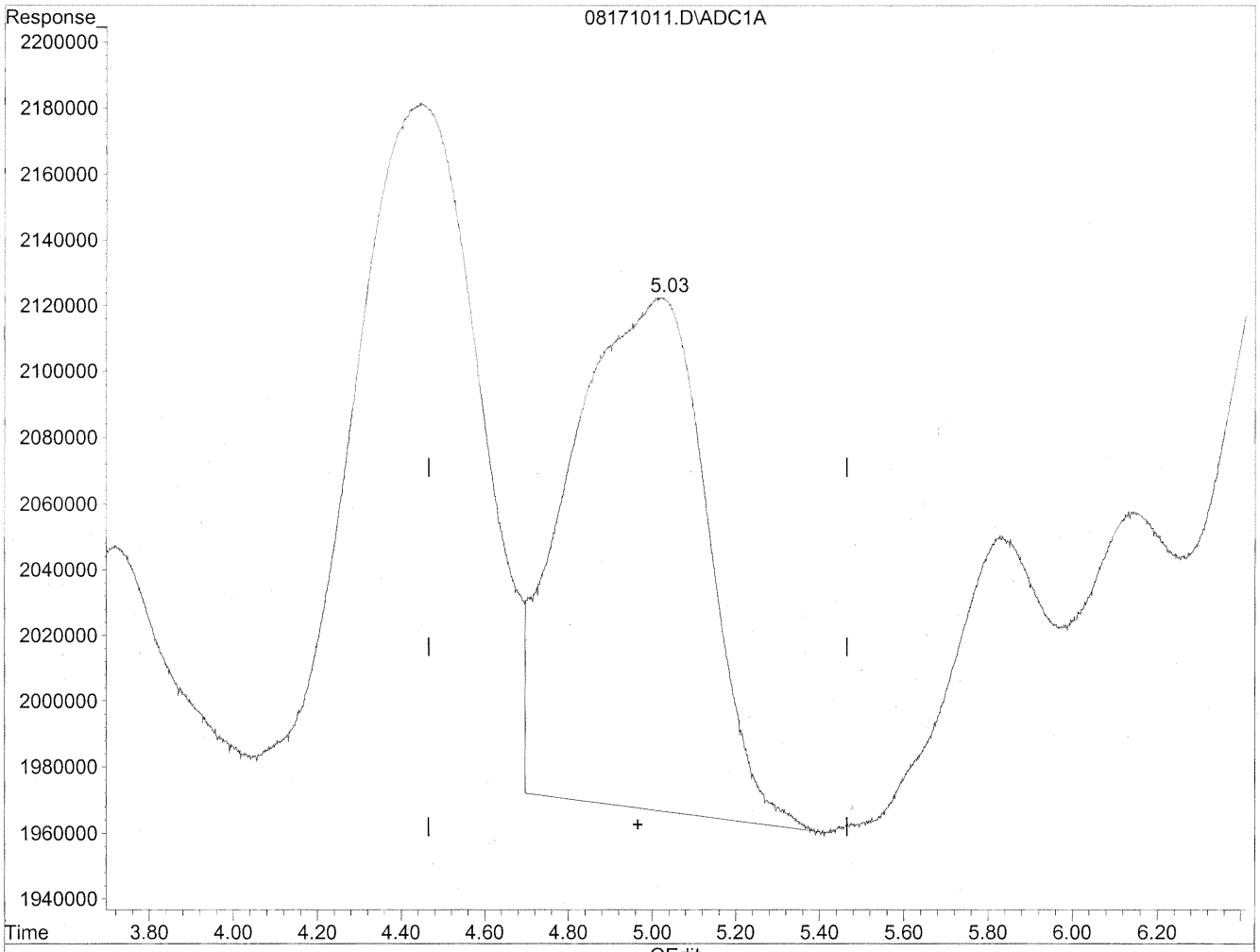
(+) = Expected Retention Time

08171011.D TO110709.M   Sat Aug 22 14:27:35 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.02min 398.543ng/ml

response 35205703

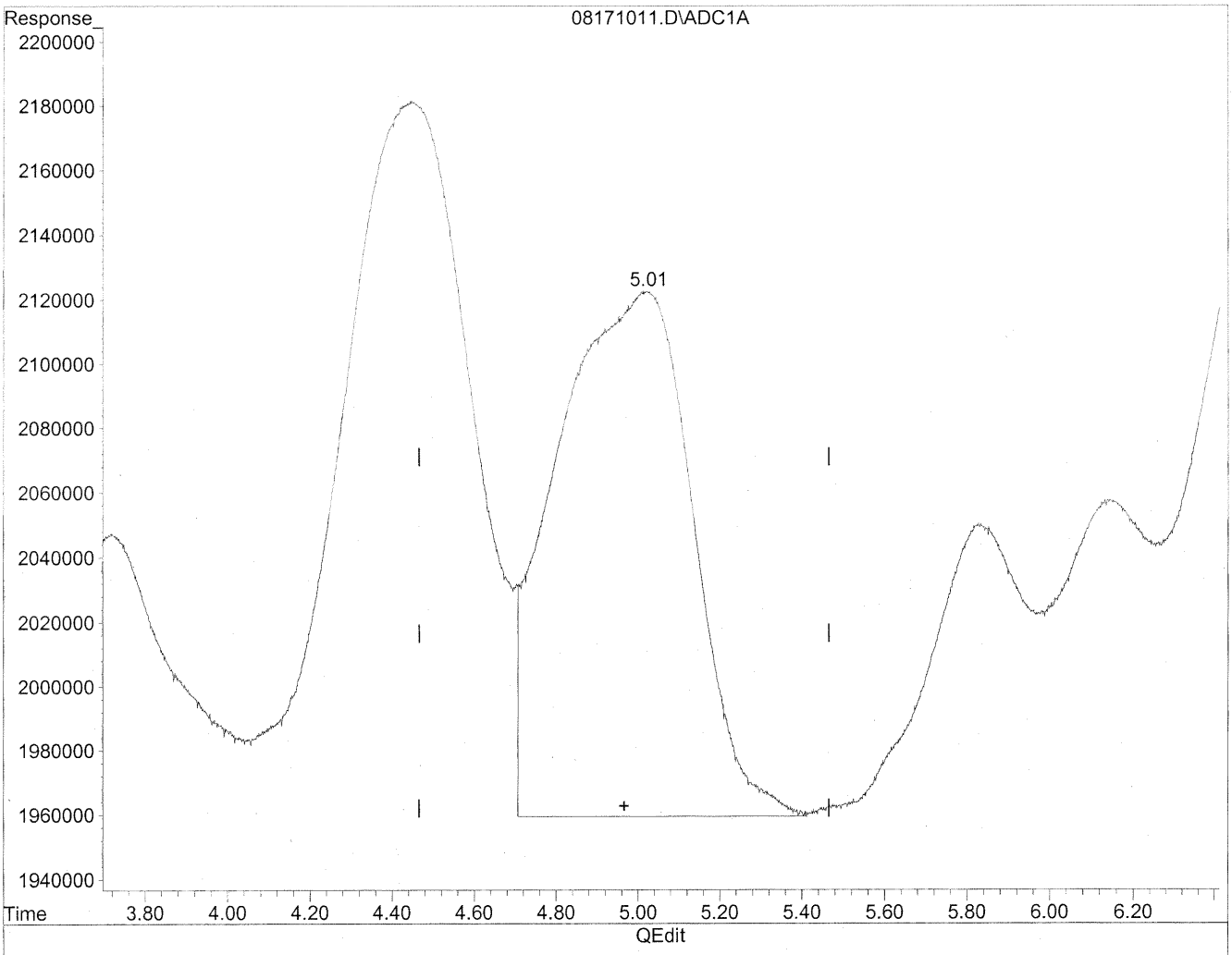
(+) = Expected Retention Time

08171011.D TO110709.M Sat Aug 22 14:27:38 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.01min 426.423ng/ml m

response 37668543

HC 8/23/09
RK 1 of 1
1 of 8/23/09

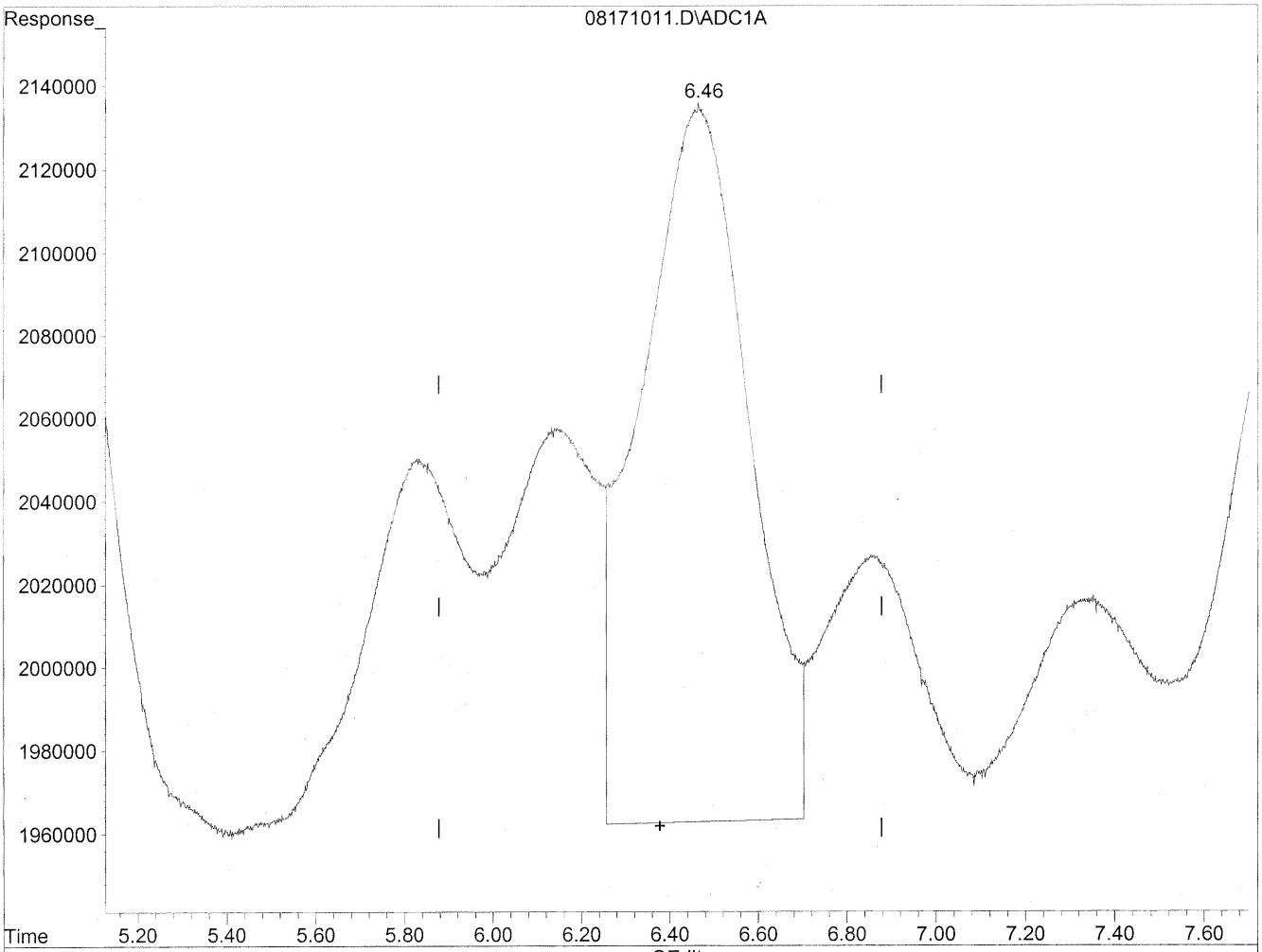
144

(+) = Expected Retention Time
08171011.D TO110709.M Sat Aug 22 14:27:51 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.46min 446.579ng/ml

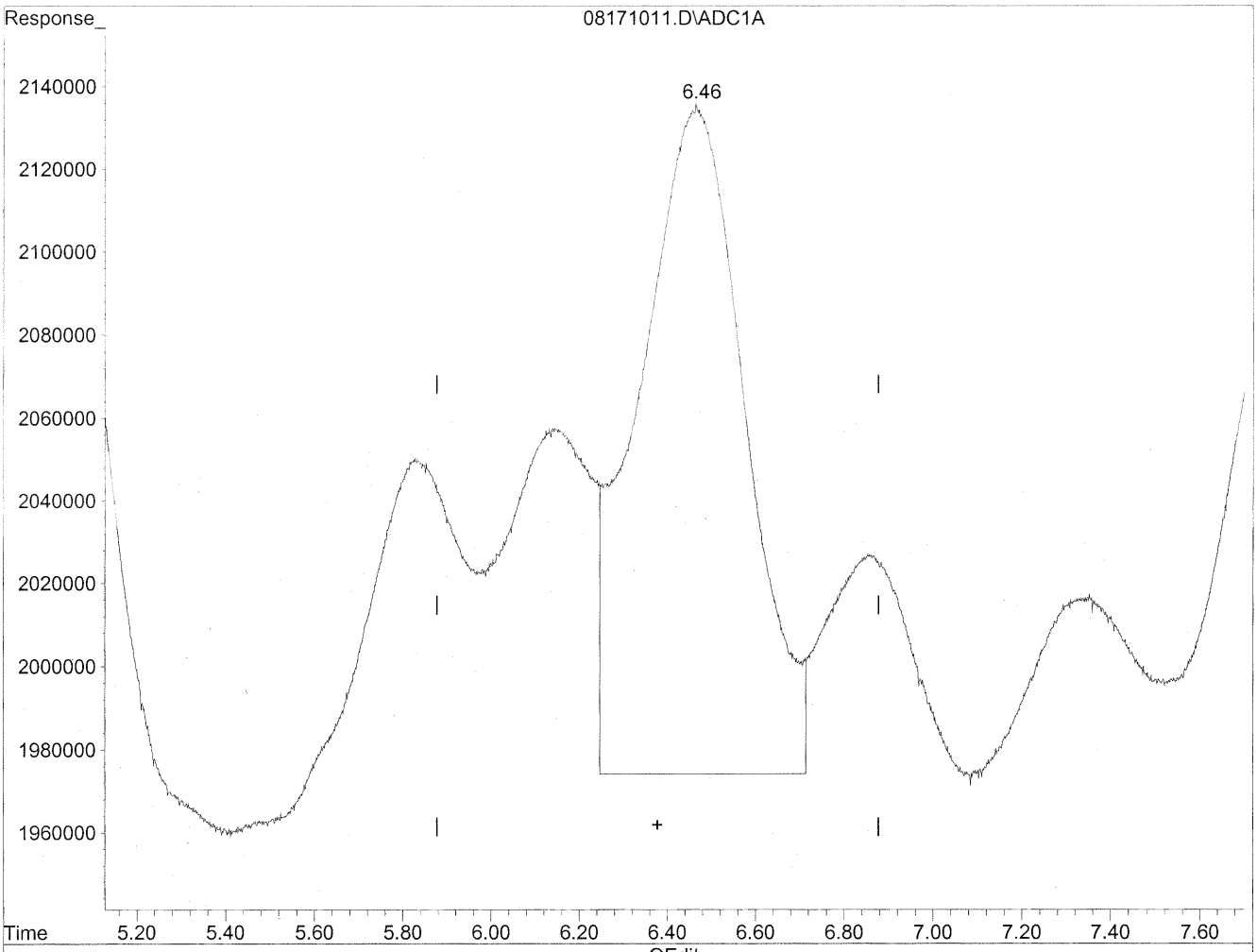
response 29415853

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.46min 409.326ng/ml m

response 26962027

HC 8/24/09
BC
KP 8/25/09

146

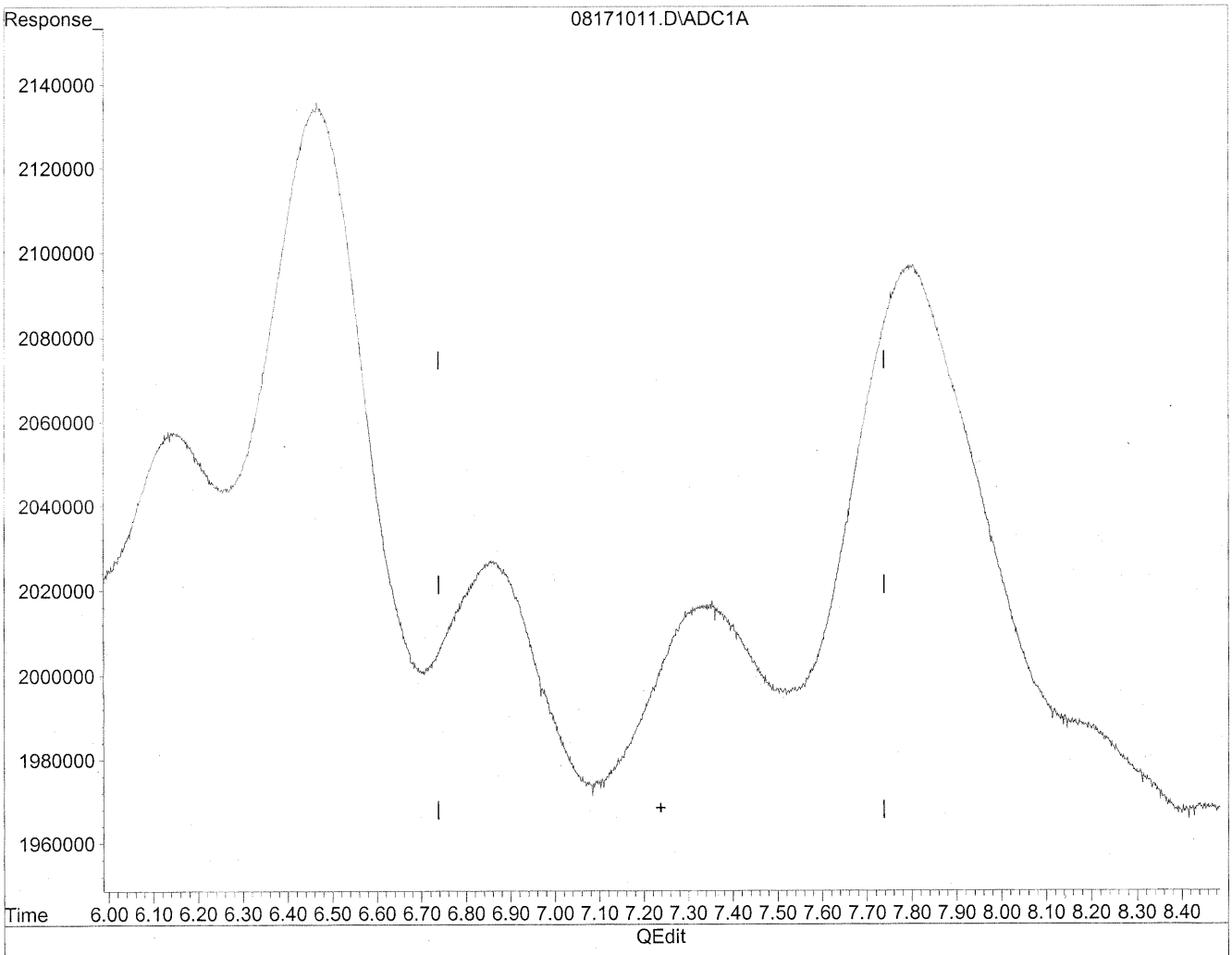
(+) = Expected Retention Time

08171011.D TO110709.M Sat Aug 22 14:28:22 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.24min 0.000ng/ml

response 0

147

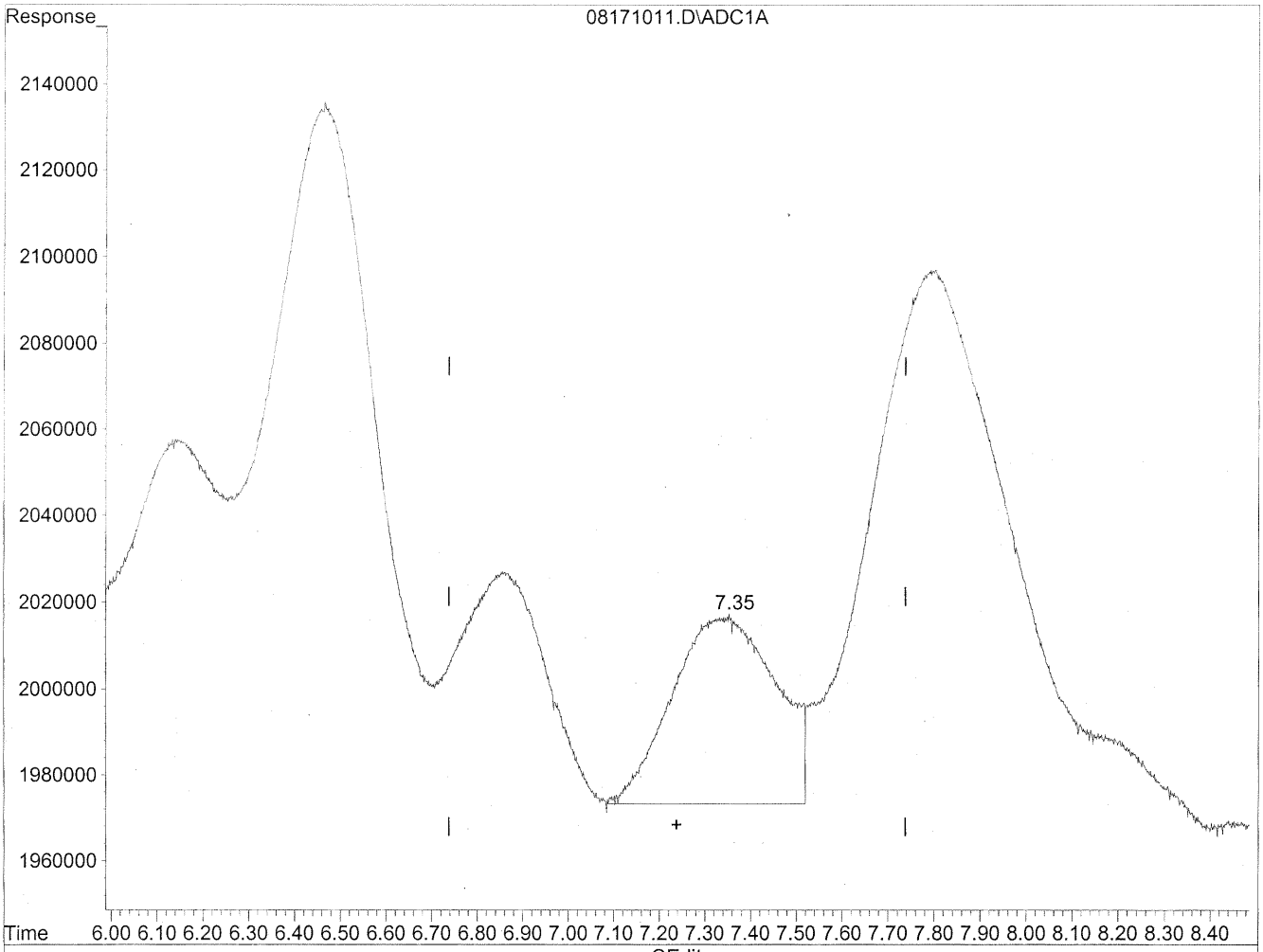
(+) = Expected Retention Time

08171011.D TO110709.M Sat Aug 22 14:28:26 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(7) Isovaleraldehyde

7.35min 88.140ng/ml m

response 6897046

HC
8/22/09
BNL
8/23/09

148

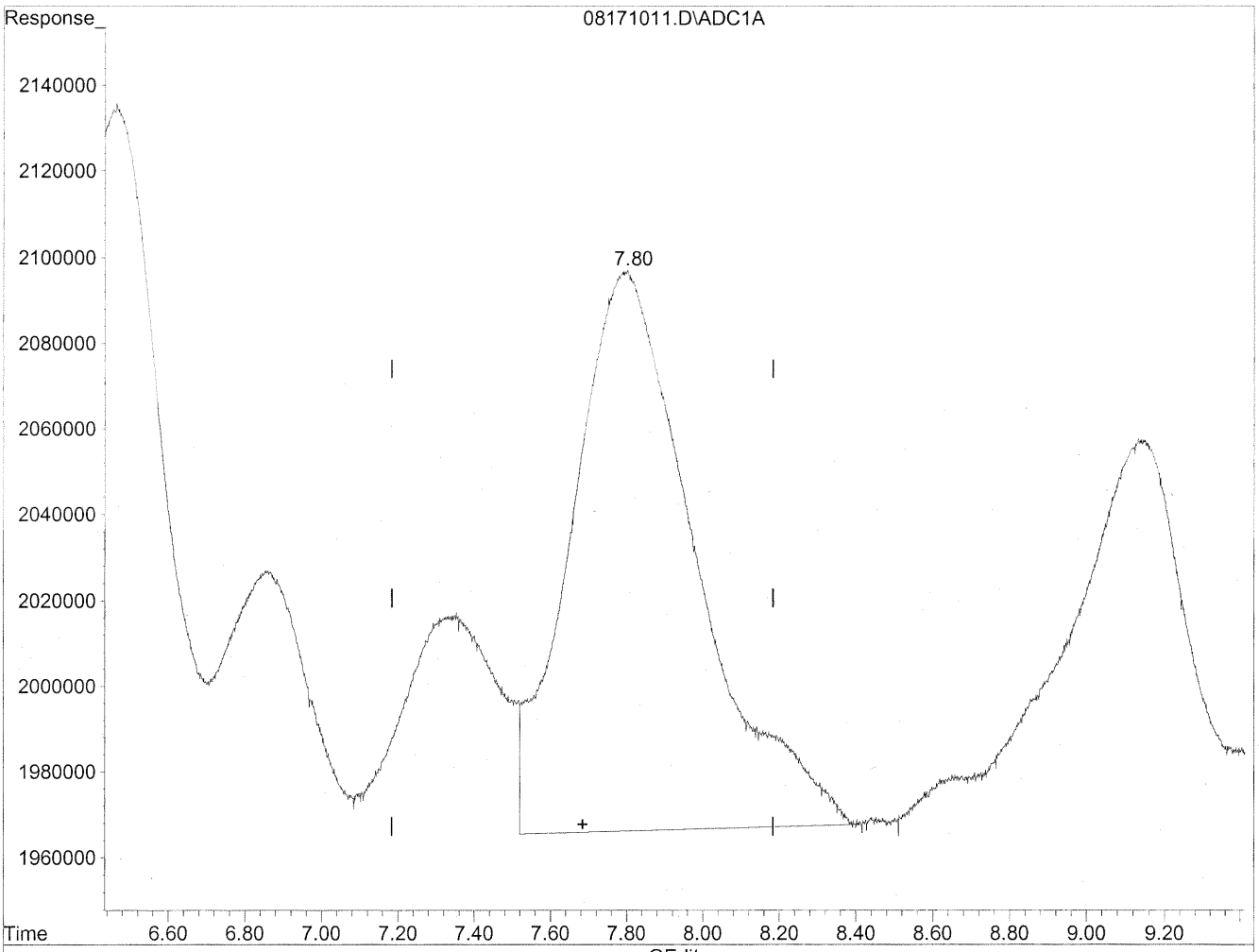
(+) = Expected Retention Time

08171011.D TO110709.M Sat Aug 22 14:28:34 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.80min 400.880ng/ml

response 29466655

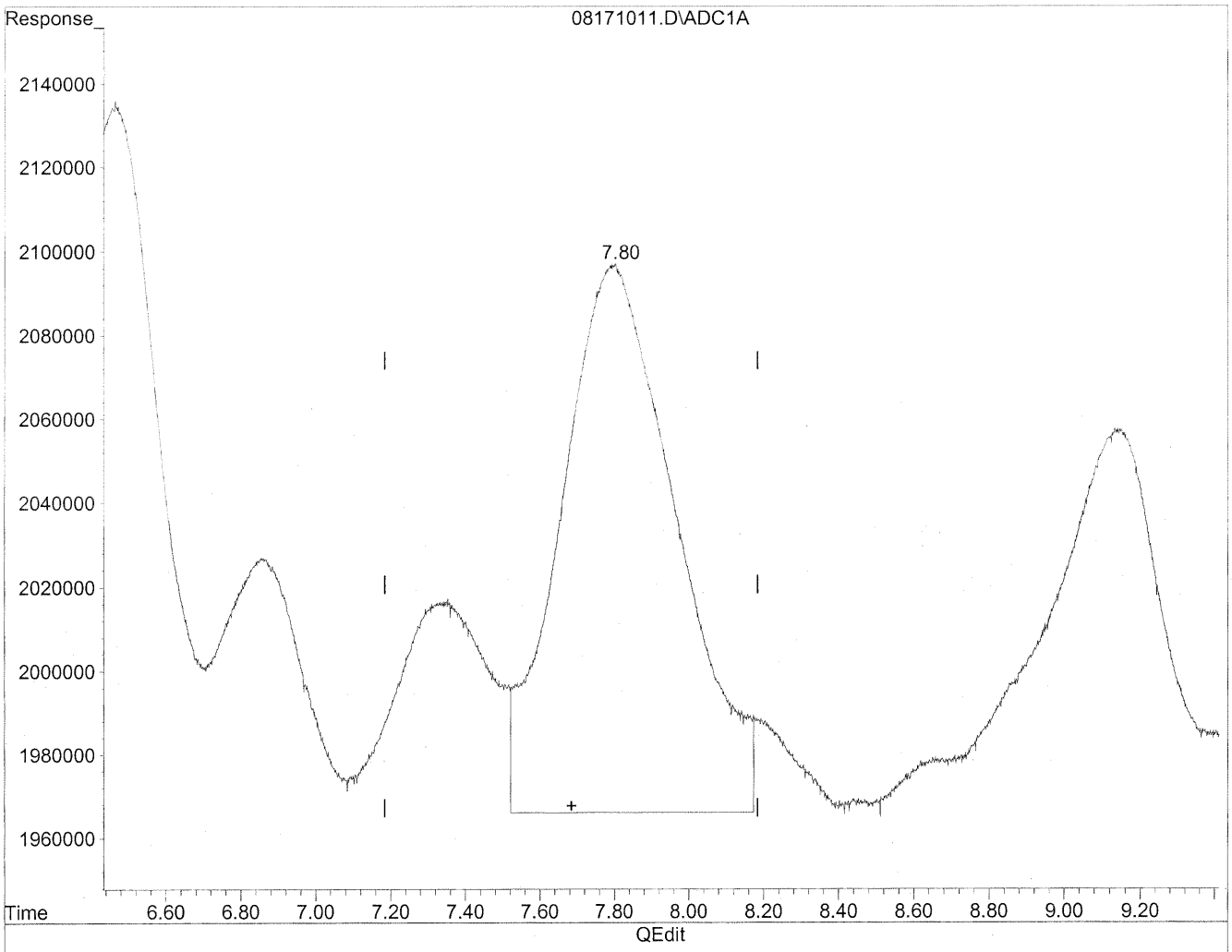
149

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.80min 380.917ng/ml m

response 27999295

HC 8/22/09 SH BC

KR 8/23/09

150

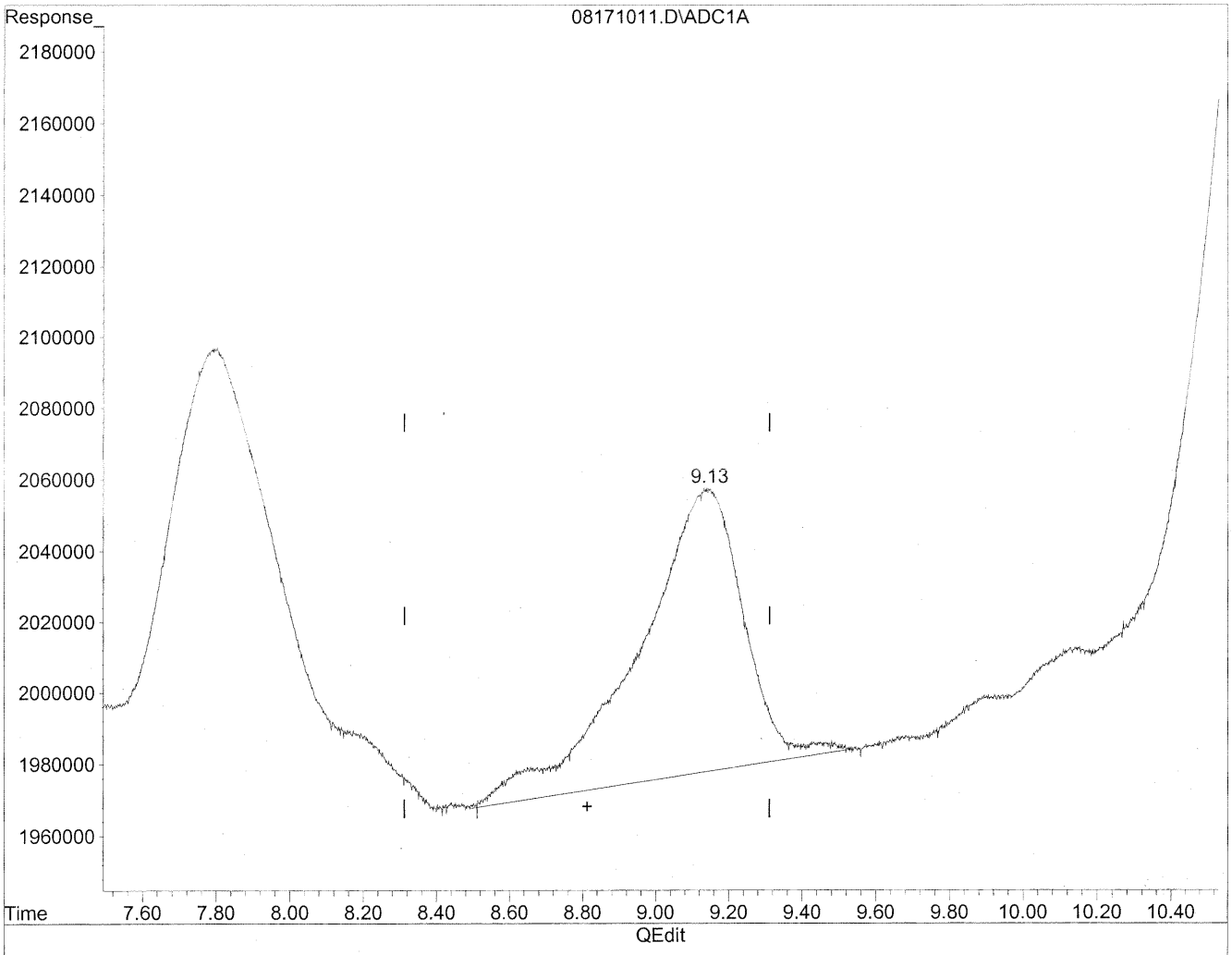
(+) = Expected Retention Time

08171011.D TO110709.M Sat Aug 22 14:28:58 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

9.14min 293.041ng/ml

response 15822881

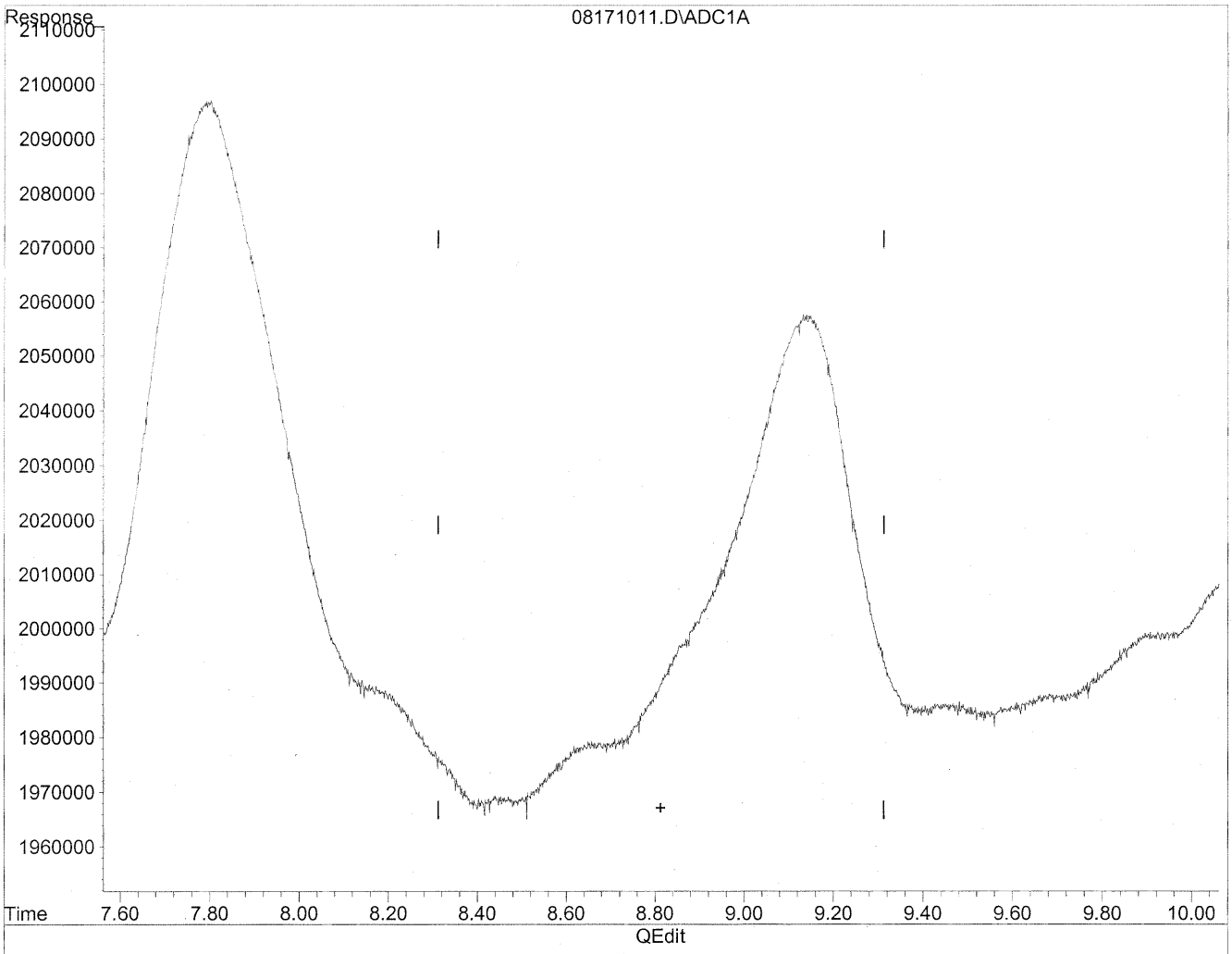
151

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09 MP

KP 8/23/09

152

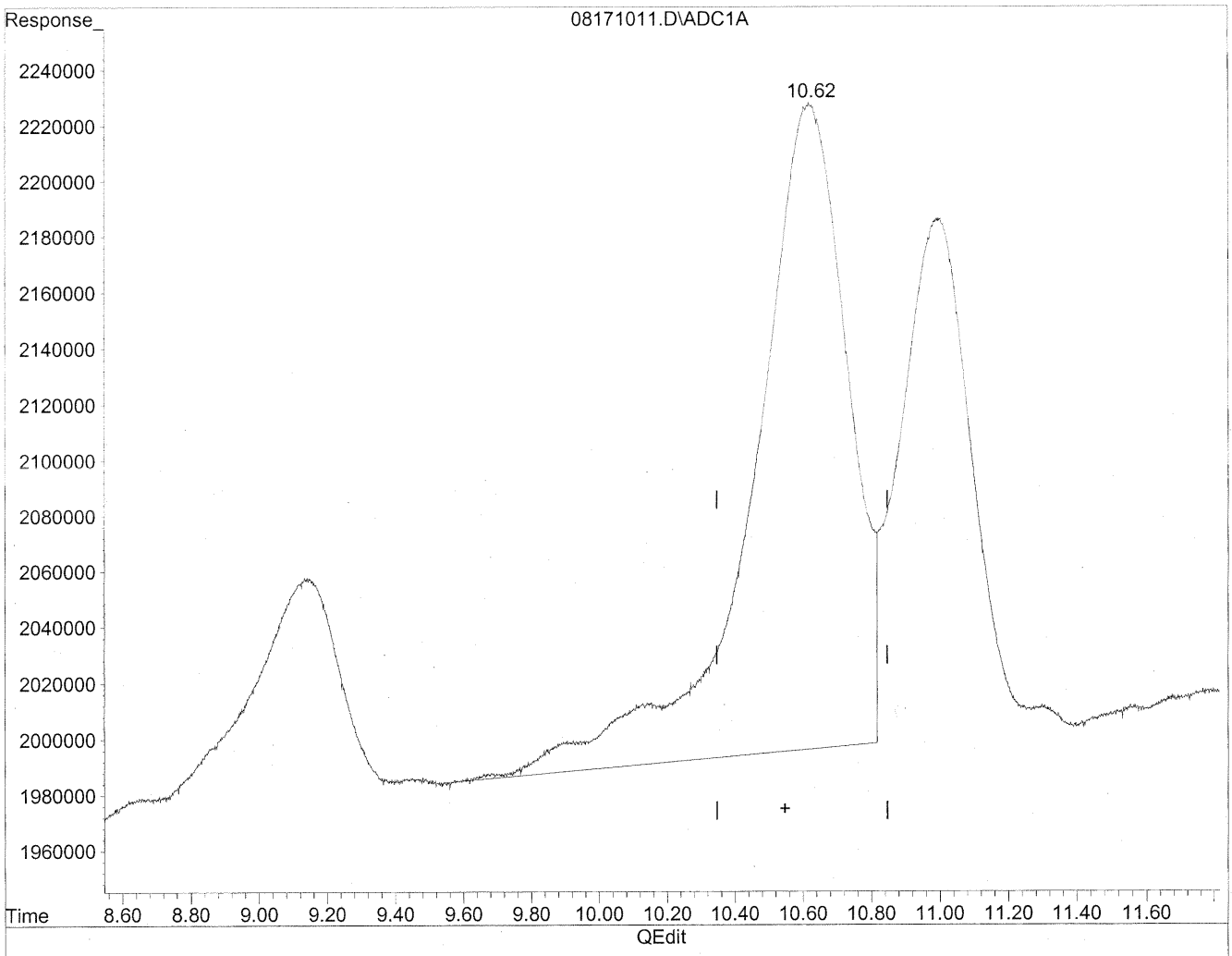
(+) = Expected Retention Time

08171011.D TO110709.M Sat Aug 22 14:29:05 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.62min 669.178ng/ml

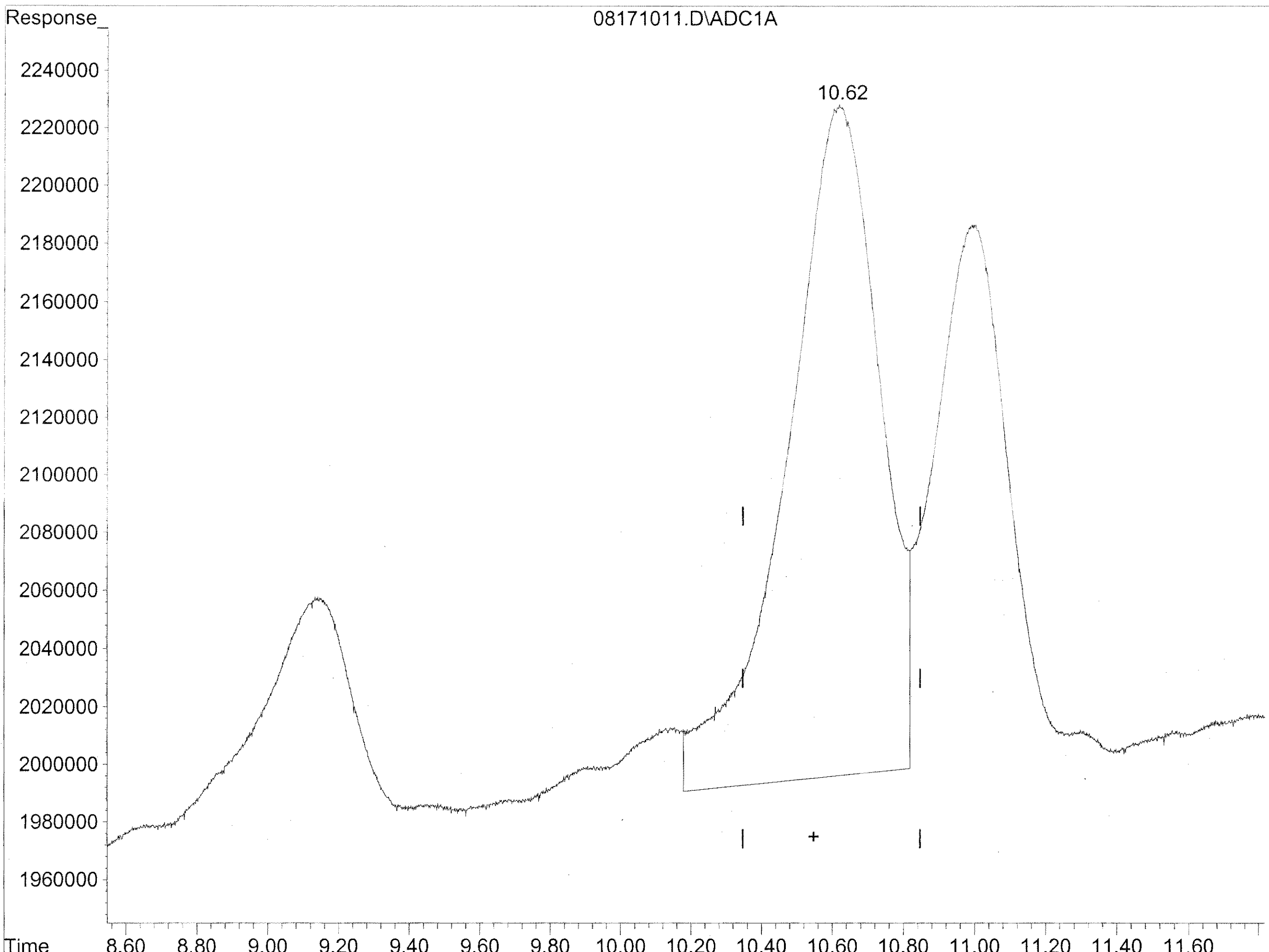
response 45064951

(+) = Expected Retention Time

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D       Vial: 24
Acq On    : 18 Aug 2009    6:24 pm                     Operator: HC
Sample    : P0902772-006 front 1.0ml                   Inst    : LC 01
Misc      :                                            Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Aug 22 14:26 19109  Quant Results File: TO110709.RES

Method        : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title         : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update   : Wed Aug 19 10:45:48 2009
Response via  : Multiple Level Calibration
```



QEdit

(11) Hexaldehyde

10.62min 624.050ng/ml m

response 42025868

HC 8/22/09 IC

KL 8/23/09

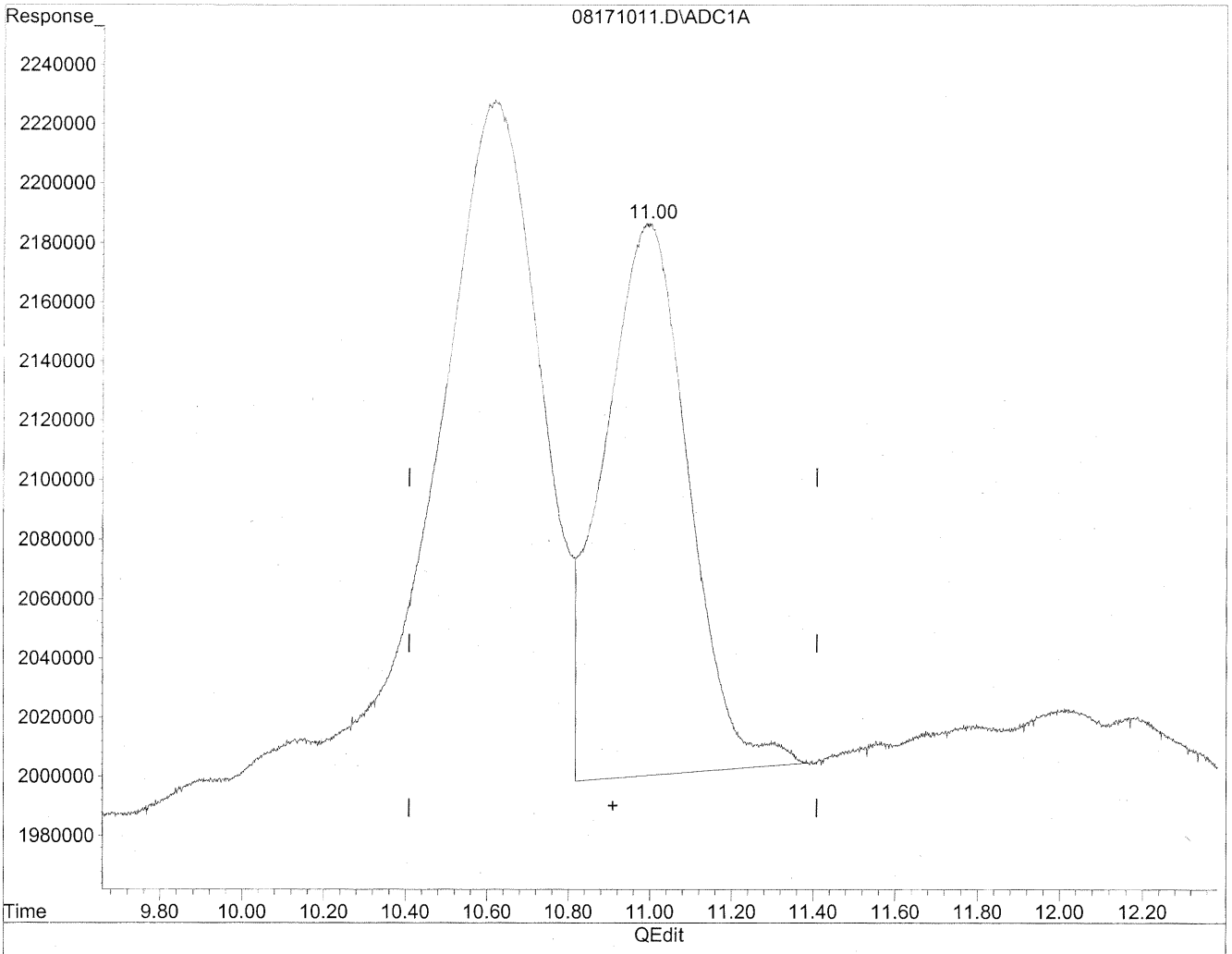
154

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.99min 549.341ng/ml

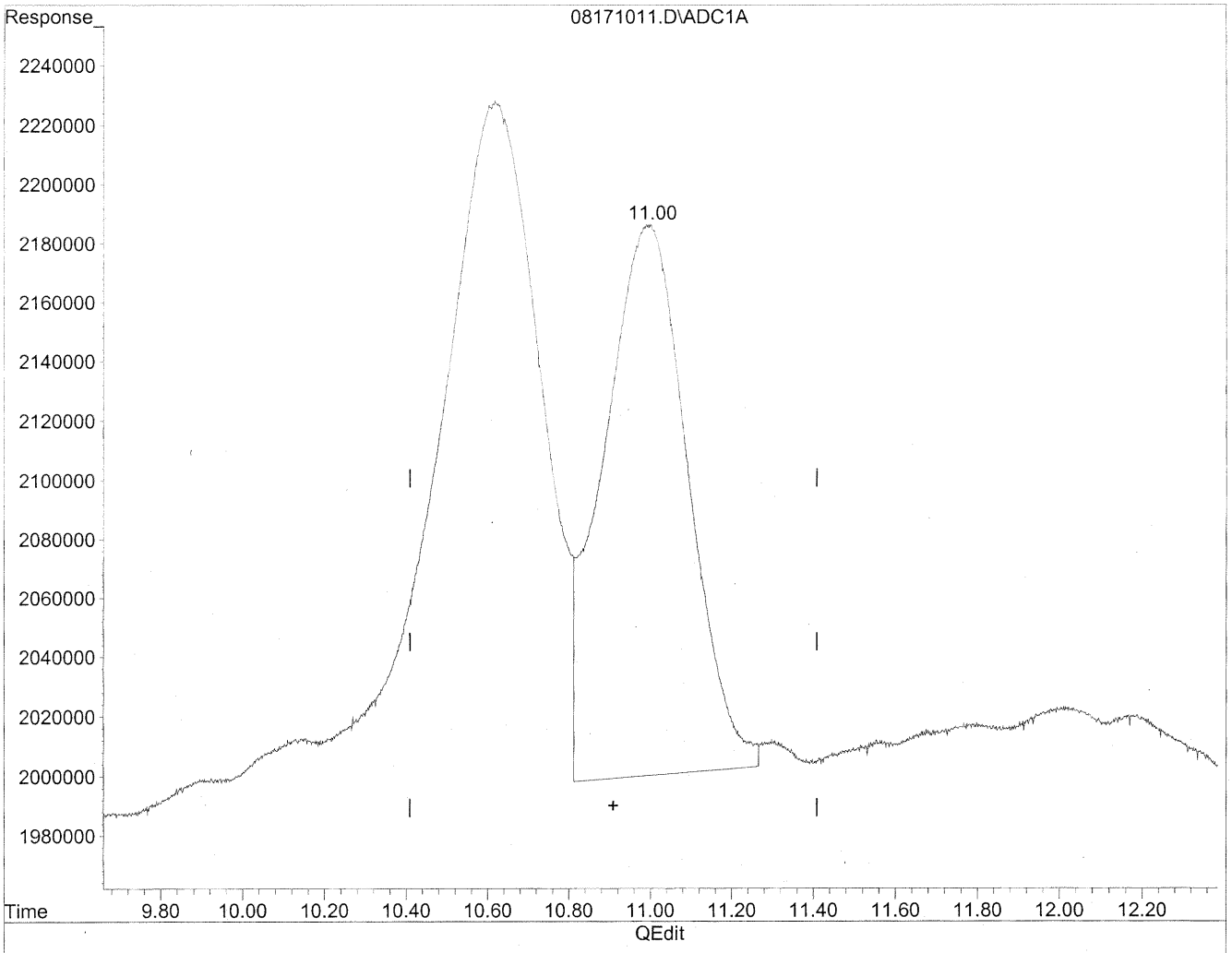
response 26925069

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171011.D Vial: 24
Acq On : 18 Aug 2009 6:24 pm Operator: HC
Sample : P0902772-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:26 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

11.00min 547.292ng/ml m

response 26824624

HC 8/22/09

KL 8/23/09

156

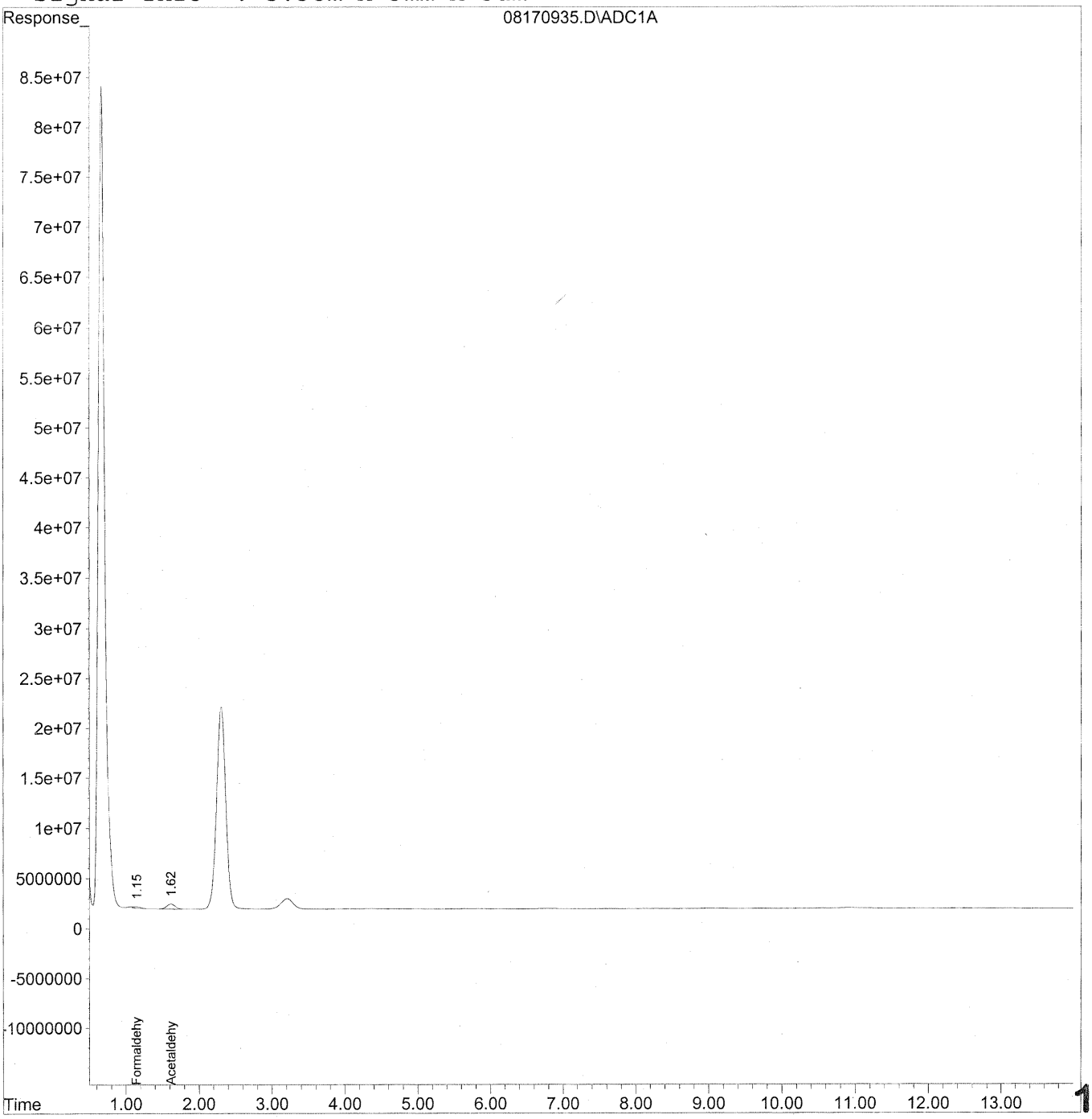
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170935.D Vial: 34
Acq On : 17 Aug 2009 11:21 pm Operator: HC
Sample : P0902772-006 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:16 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



157

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170935.D Vial: 34
Acq On : 17 Aug 2009 11:21 pm Operator: HC
Sample : P0902772-006 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:16 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

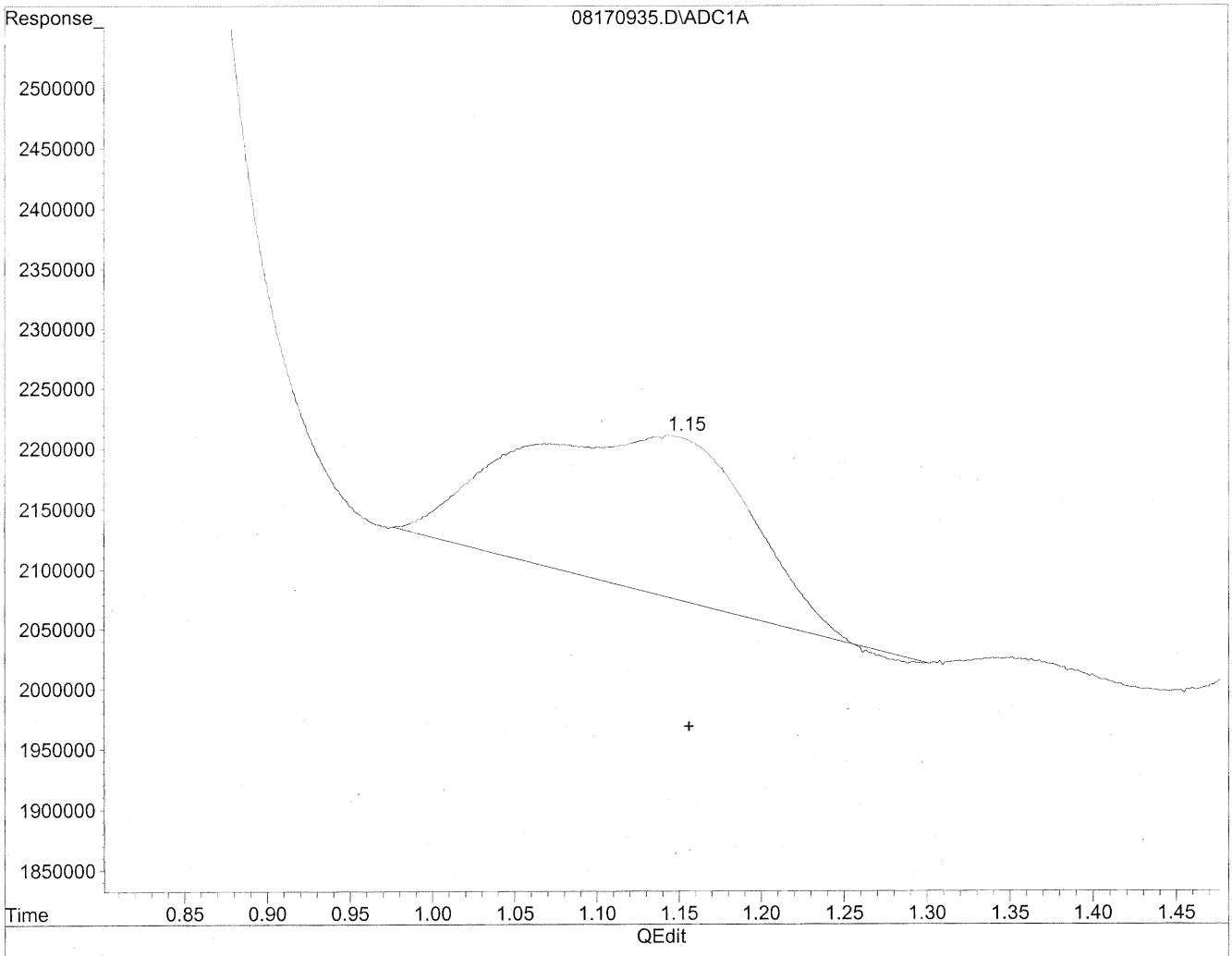
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.15	9298193	50.649	ng/mlm
2)	Acetaldehyde	1.62	39812800	283.924	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170935.D Vial: 34
Acq On : 17 Aug 2009 11:21 pm Operator: HC
Sample : P0902772-006 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:15 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.14min 70.535ng/ml

response 12948864

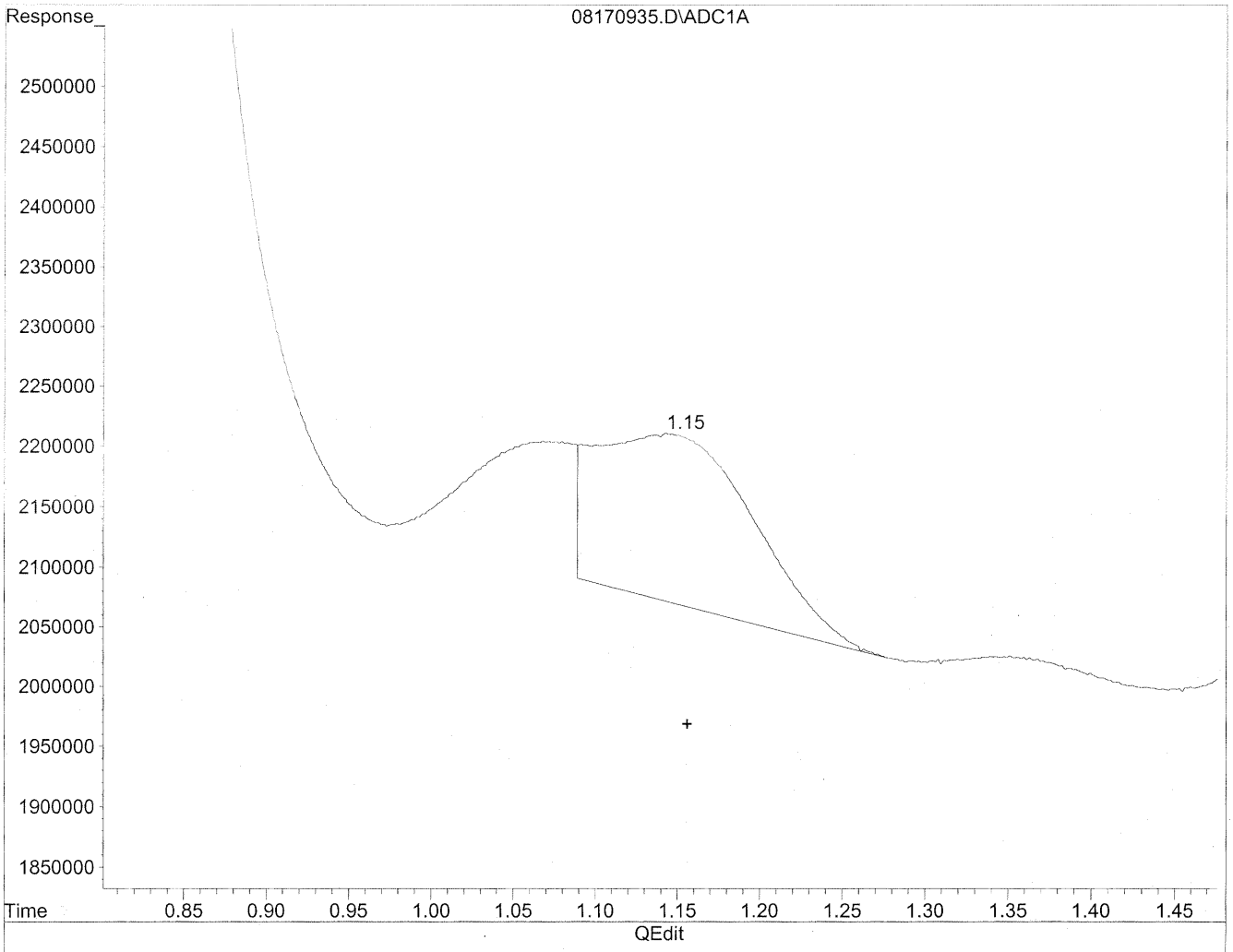
159

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170935.D Vial: 34
Acq On : 17 Aug 2009 11:21 pm Operator: HC
Sample : P0902772-006 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:15 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.15min 50.649ng/ml m

response 9298193

HC 8/21/09
SP
KK 8/23/09

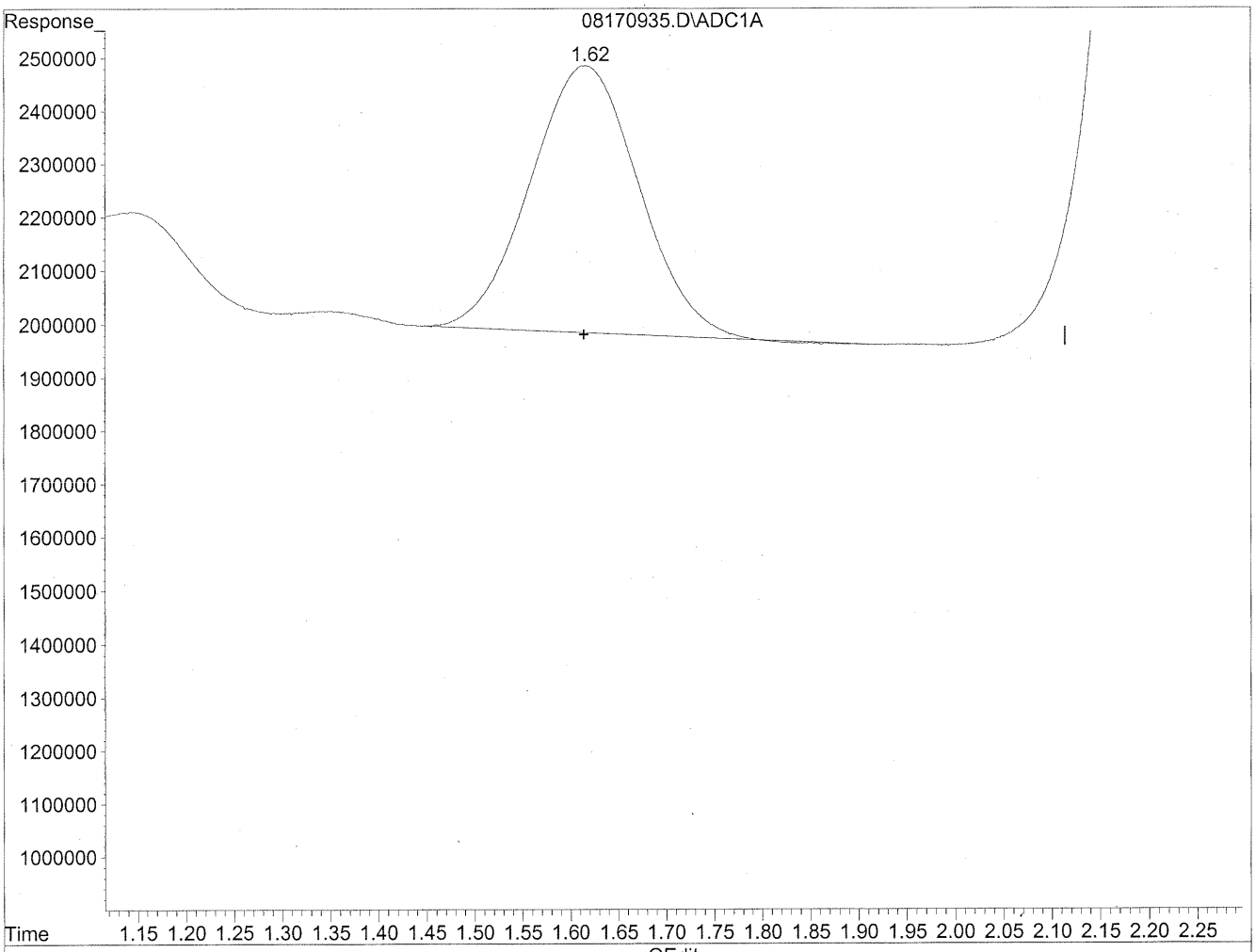
(+) = Expected Retention Time

08170935.D TO110709.M Fri Aug 21 18:16:14 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170935.D Vial: 34
Acq On : 17 Aug 2009 11:21 pm Operator: HC
Sample : P0902772-006 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:15 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(2) Acetaldehyde

1.62min 280.788ng/ml

response 39373120

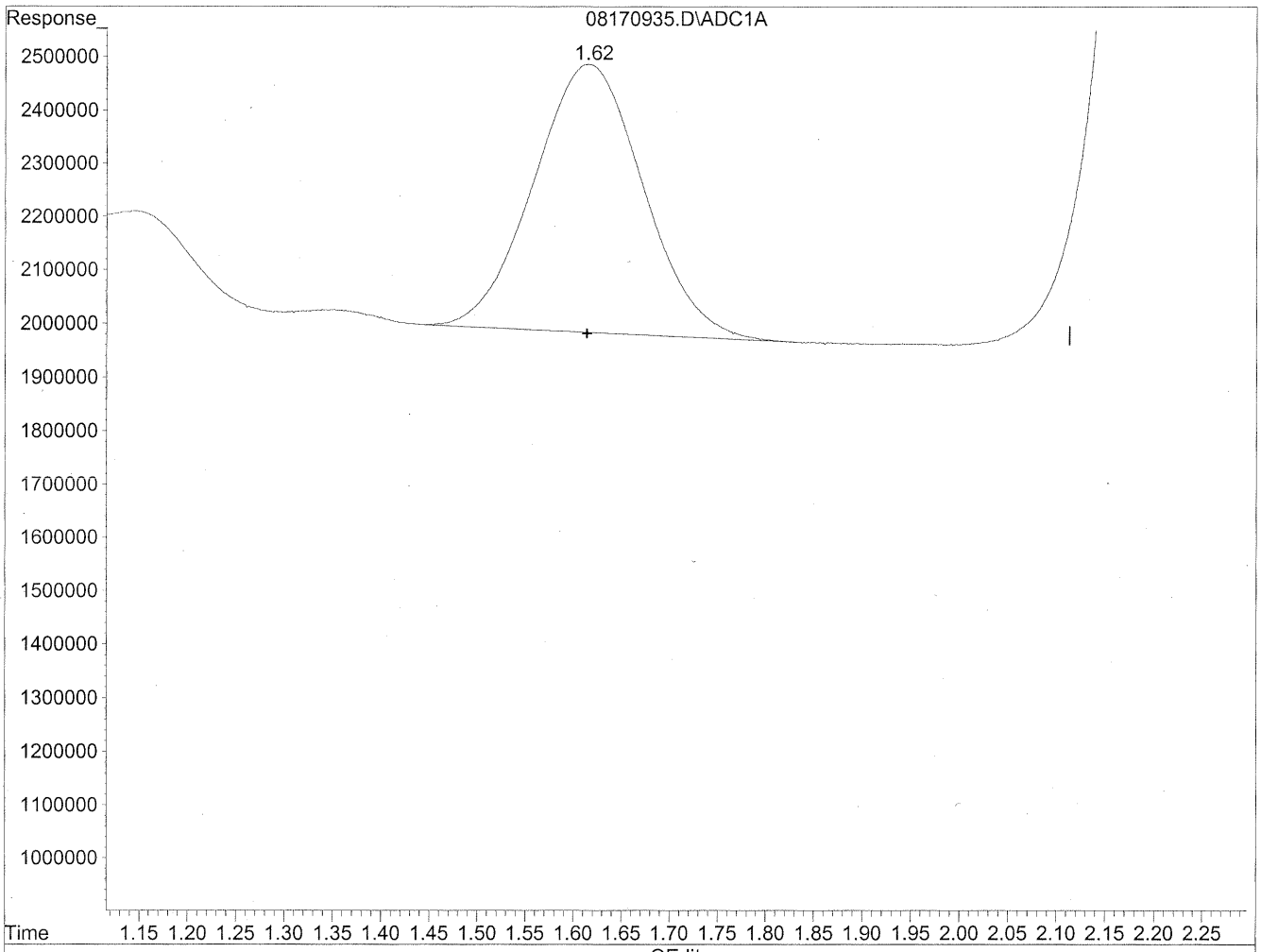
(+) = Expected Retention Time

08170935.D TO110709.M Fri Aug 21 18:16:20 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170935.D Vial: 34
Acq On : 17 Aug 2009 11:21 pm Operator: HC
Sample : P0902772-006 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:15 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(2) Acetaldehyde

1.62min 283.924ng/ml m

response 39812800

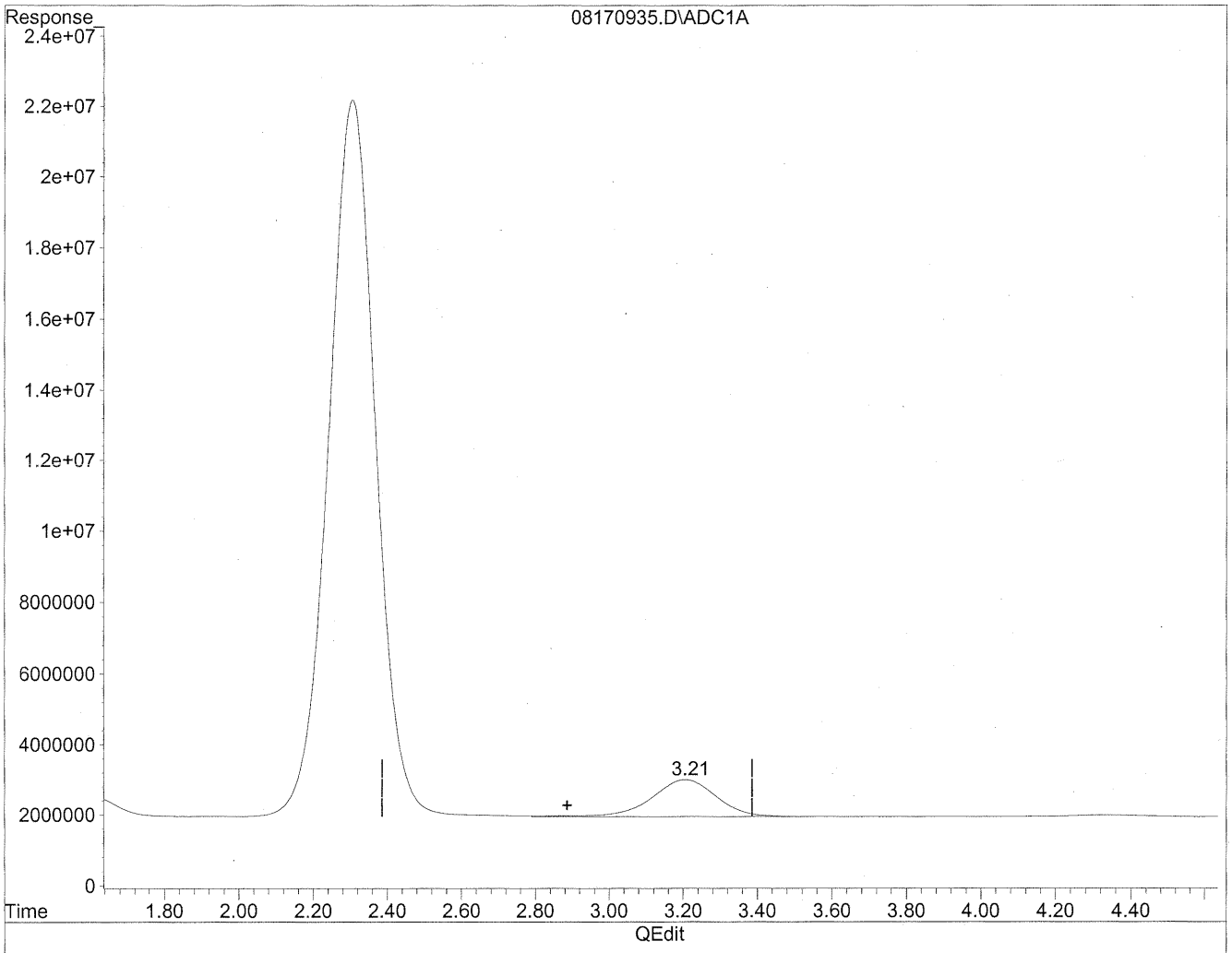
(+) = Expected Retention Time

08170935.D TO110709.M Fri Aug 21 18:16:31 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170935.D Vial: 34
Acq On : 17 Aug 2009 11:21 pm Operator: HC
Sample : P0902772-006 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:15 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.21min 1191.552ng/ml

response 127132890

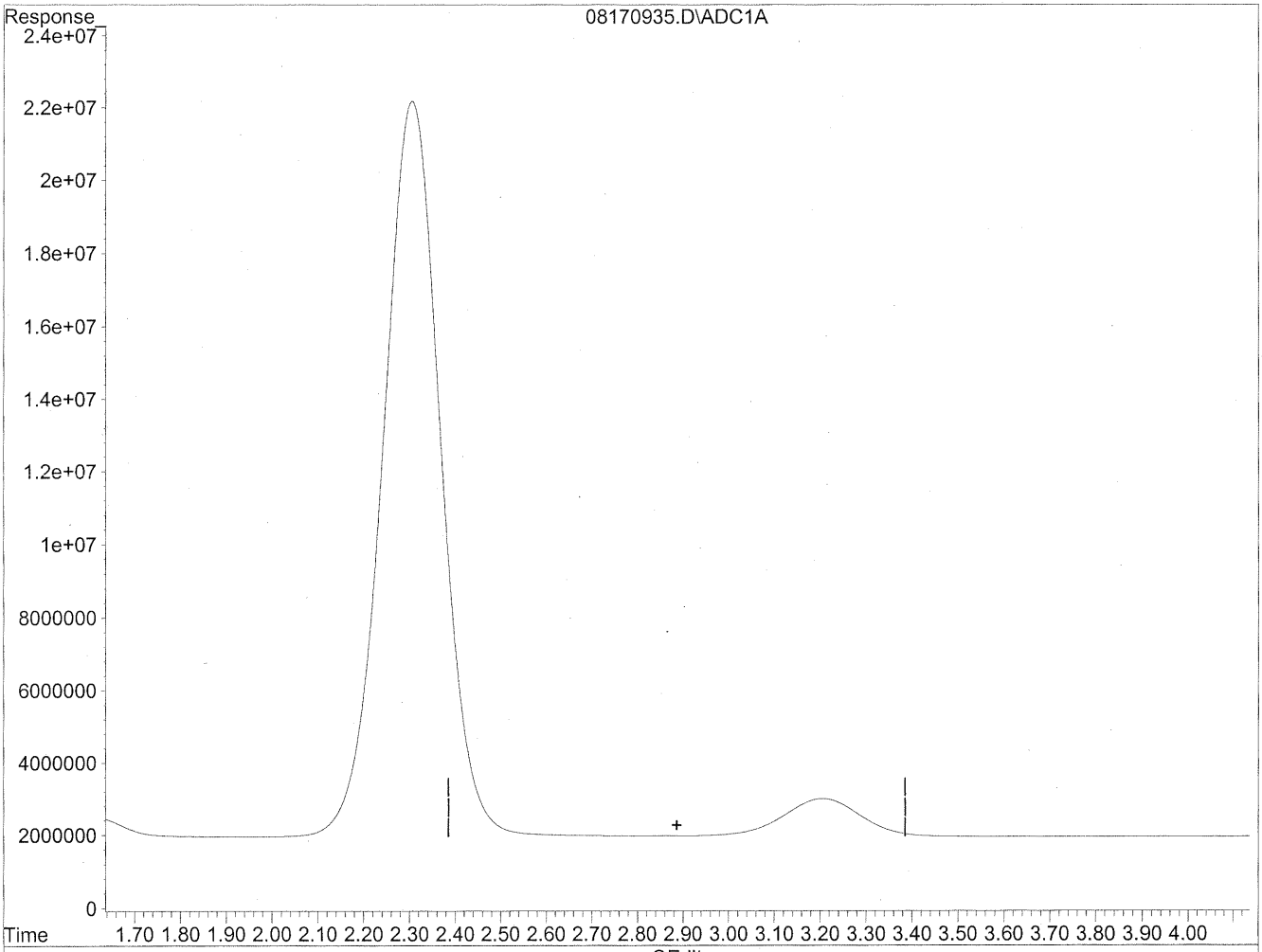
(+) = Expected Retention Time

08170935.D TO110709.M Fri Aug 21 18:16:34 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170935.D Vial: 34
Acq On : 17 Aug 2009 11:21 pm Operator: HC
Sample : P0902772-006 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:15 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/21/09 MP

KP 8/23/09

164

(+) = Expected Retention Time

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **100443**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P0902772-007

Test Code: EPA Method TO-11A
Instrument ID: Waters LC Module I Plus/UV_Vis 360/LC1
Analyst: Hani Cherazaie
Sampling Media: Silica Gel DNPH Tube
Test Notes: **BC**

Date Collected: 8/10/09
Date Received: 8/12/09
Date Analyzed: 8/18 - 8/19/09
Desorption Volume: 1.0 ml
Volume Sampled: 106.3 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**12,000**	**110**	0.94	**93**	0.77	
75-07-0	**Acetaldehyde**	**15,000**	**140**	0.94	**77**	0.52	**BT**
123-38-6	**Propionaldehyde**	**880**	**8.2**	0.94	**3.5**	0.40	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.94	ND	0.33	
123-72-8	**Butyraldehyde**	**530**	**5.0**	0.94	**1.7**	0.32	
100-52-7	**Benzaldehyde**	**1,000**	**9.8**	0.94	**2.3**	0.22	
590-86-3	**Isovaleraldehyde**	**280**	**2.6**	0.94	**0.74**	0.27	
110-62-3	**Valeraldehyde**	**1,100**	**9.9**	0.94	**2.8**	0.27	
529-20-4	o-Tolualdehyde	< 100	ND	0.94	ND	0.19	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	1.9	ND	0.38	
66-25-1	**n-Hexaldehyde**	**2,700**	**25**	0.94	**6.1**	0.23	
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	ND	0.94	ND	0.17	

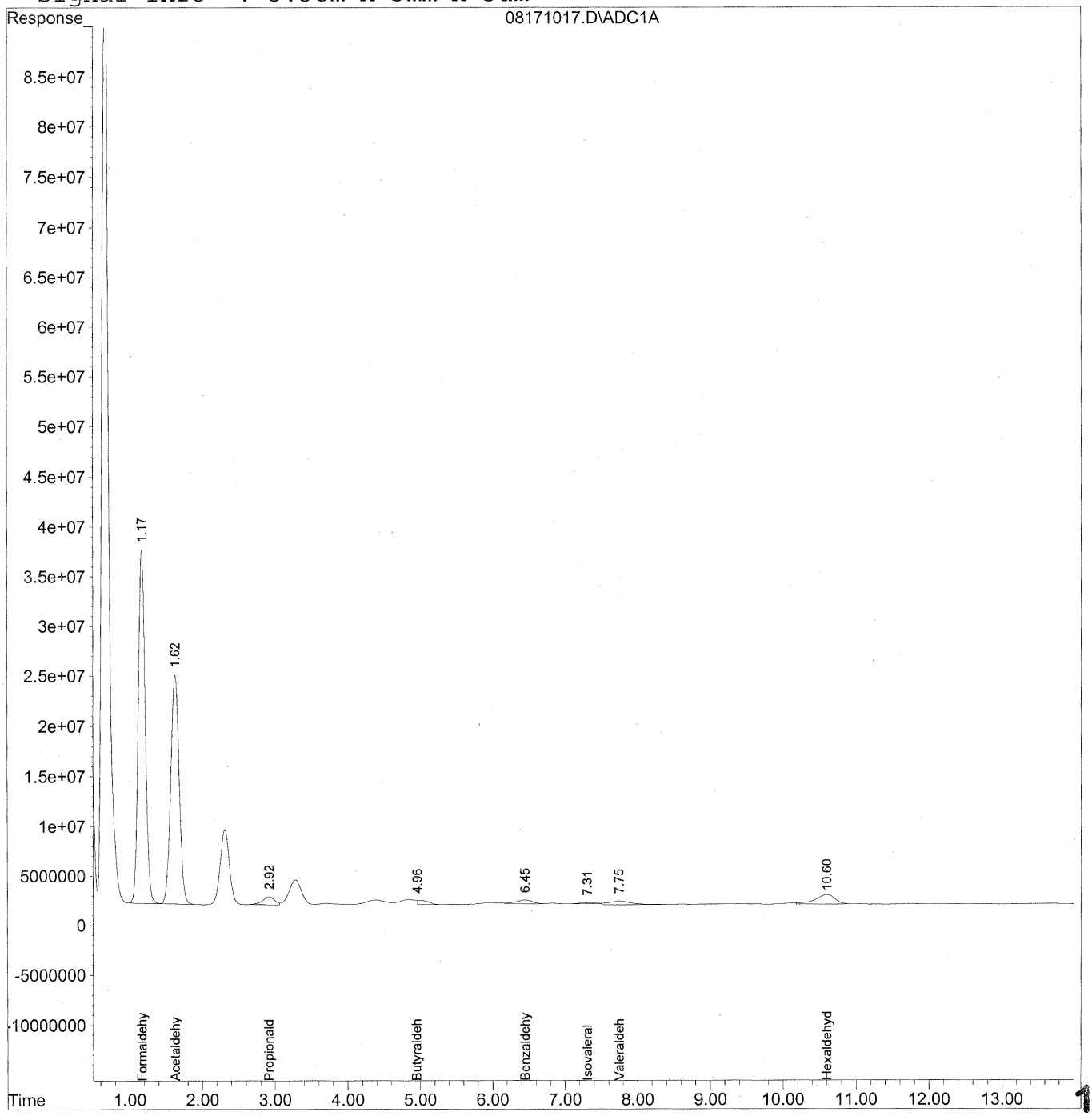
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
BT = Results indicated possible breakthrough; back section > 10% front section.

Verified By: _____ Date: 8/25/09

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:32 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



166

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:32 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

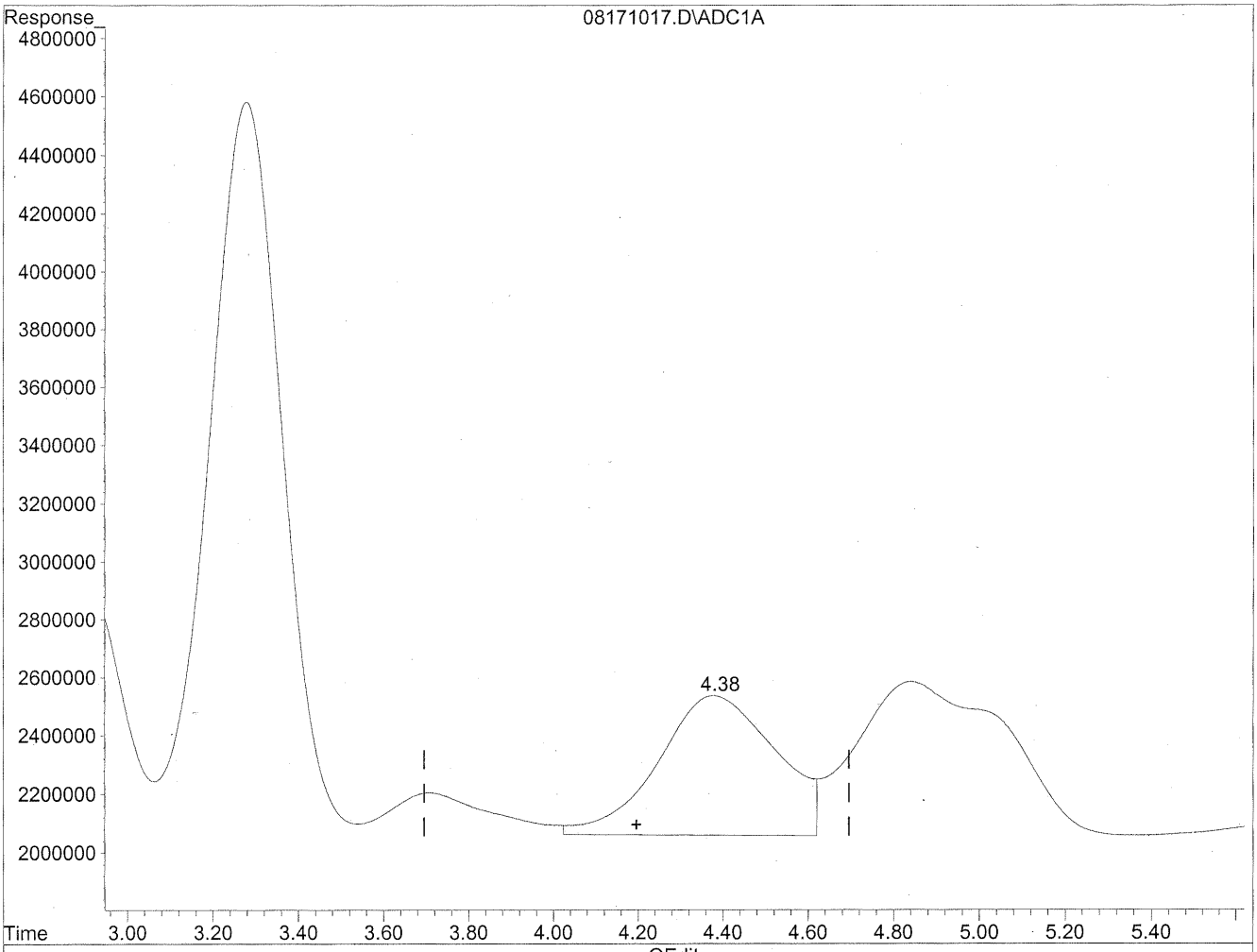
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	2332476913	12705.410	ng/ml
2)	Acetaldehyde	1.62	1842343393	13138.620	ng/ml
3)	Propionaldehyde	2.91	93536572	876.671	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	4.96	46593661	527.459	ng/mlm
6)	Benzaldehyde	6.45	68431305	1038.895	ng/mlm
7)	Isovaleraldehyde	7.31	21716251	277.520	ng/mlm
8)	Valeraldehyde	7.75	77207887	1050.376	ng/mlm
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	10.60	178815052	2655.258	ng/mlm
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/mld

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.38min 942.021ng/ml

response 91767146

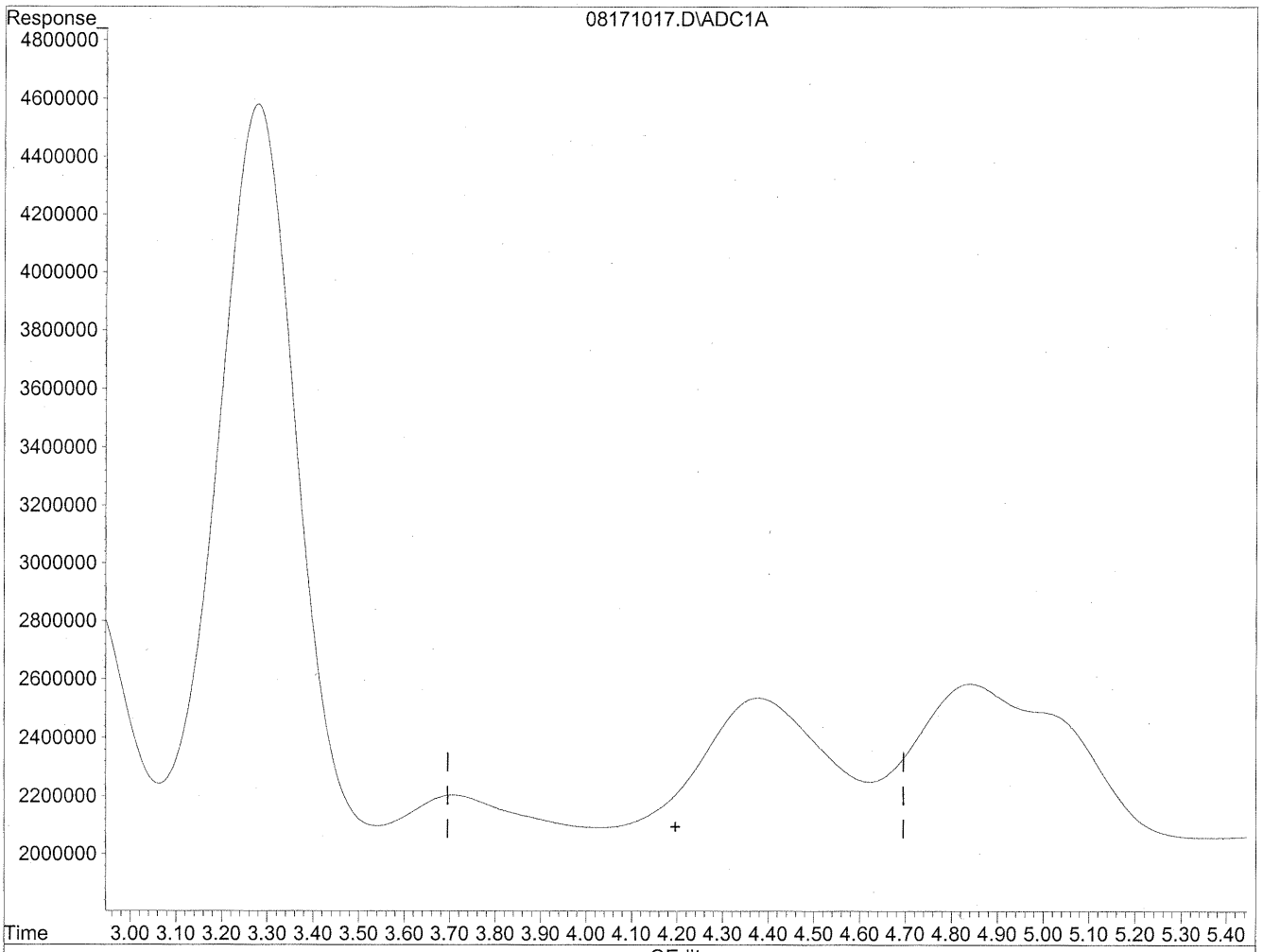
(+) = Expected Retention Time

08171017.D TO110709.M Sat Aug 22 14:31:26 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09
MP
8/23/09

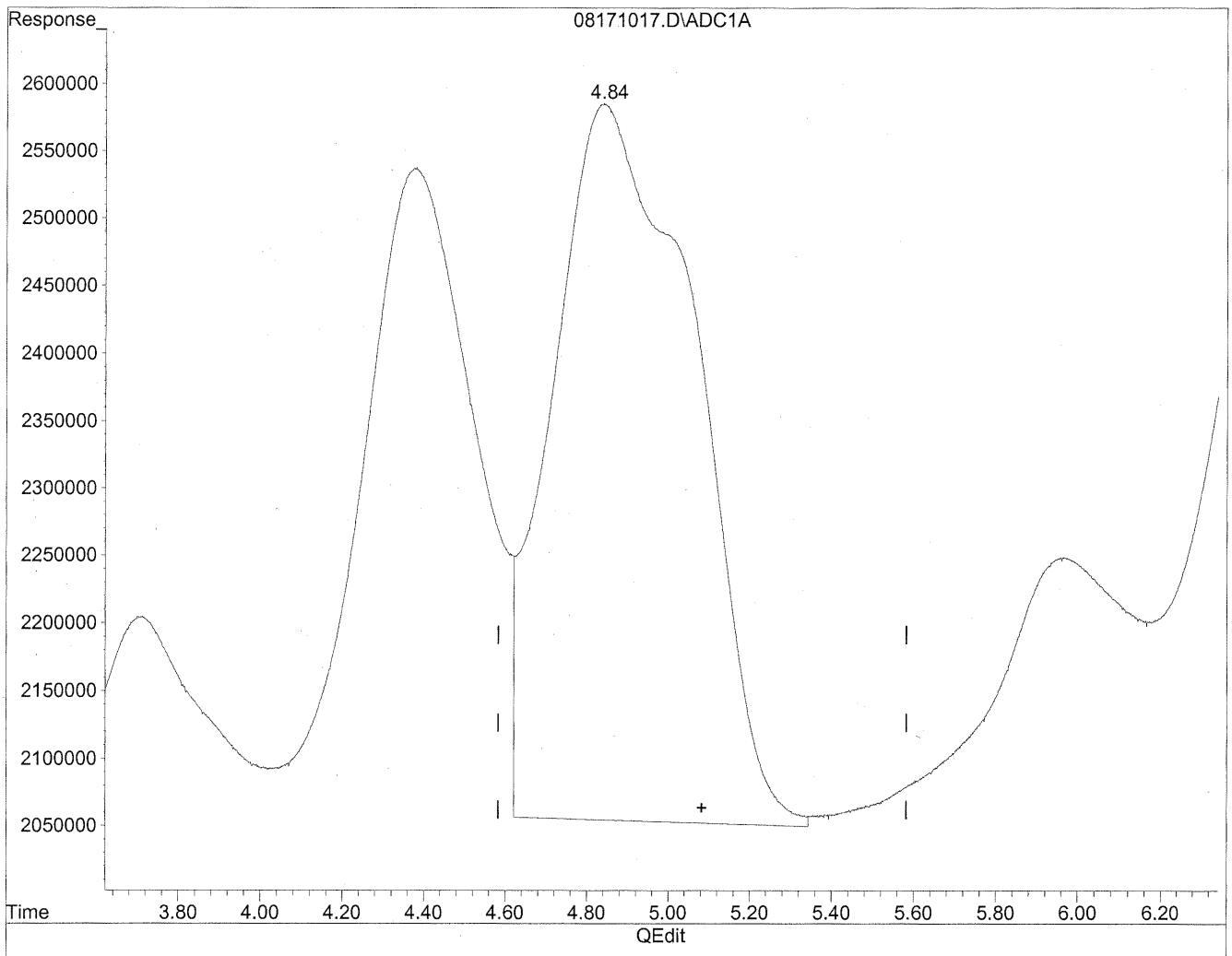
169

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.84min 1461.762ng/ml

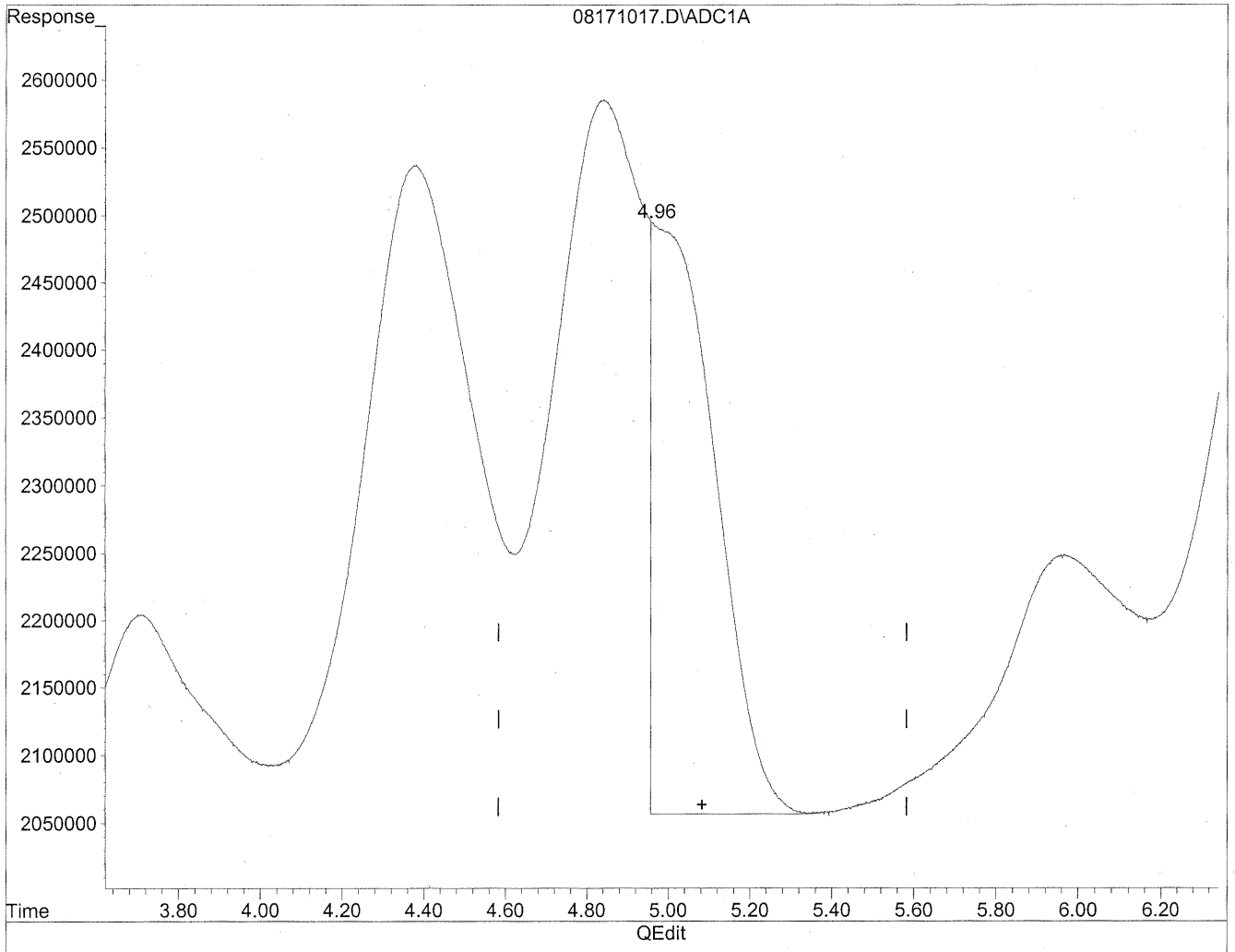
response 129126401

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.96min 527.459ng/ml m

response 46593661

tlc 8/22/09
sp
KB 8/23/09

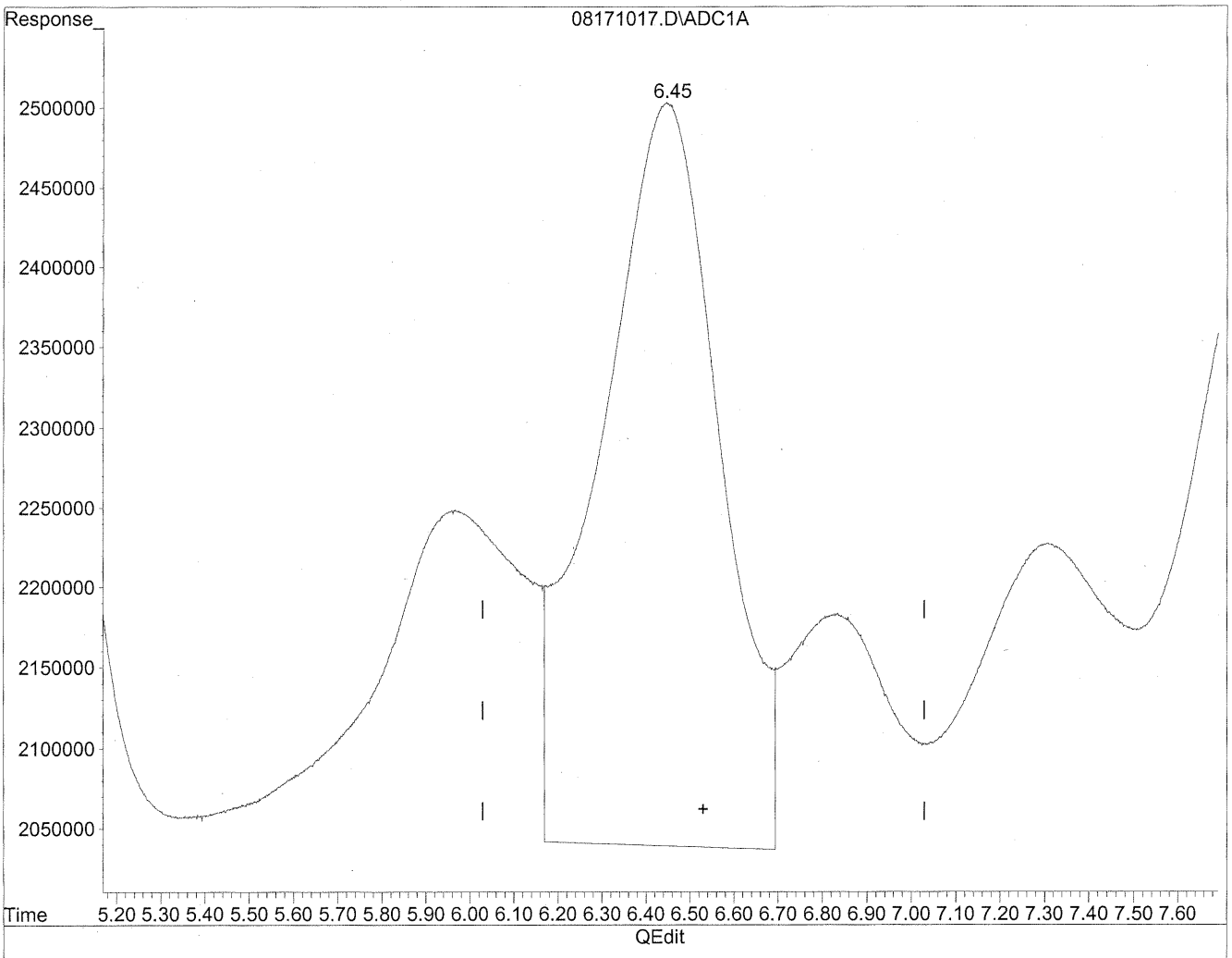
171

(+) = Expected Retention Time
08171017.D TO110709.M Sat Aug 22 14:31:41 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.45min 1330.285ng/ml

response 87624977

172

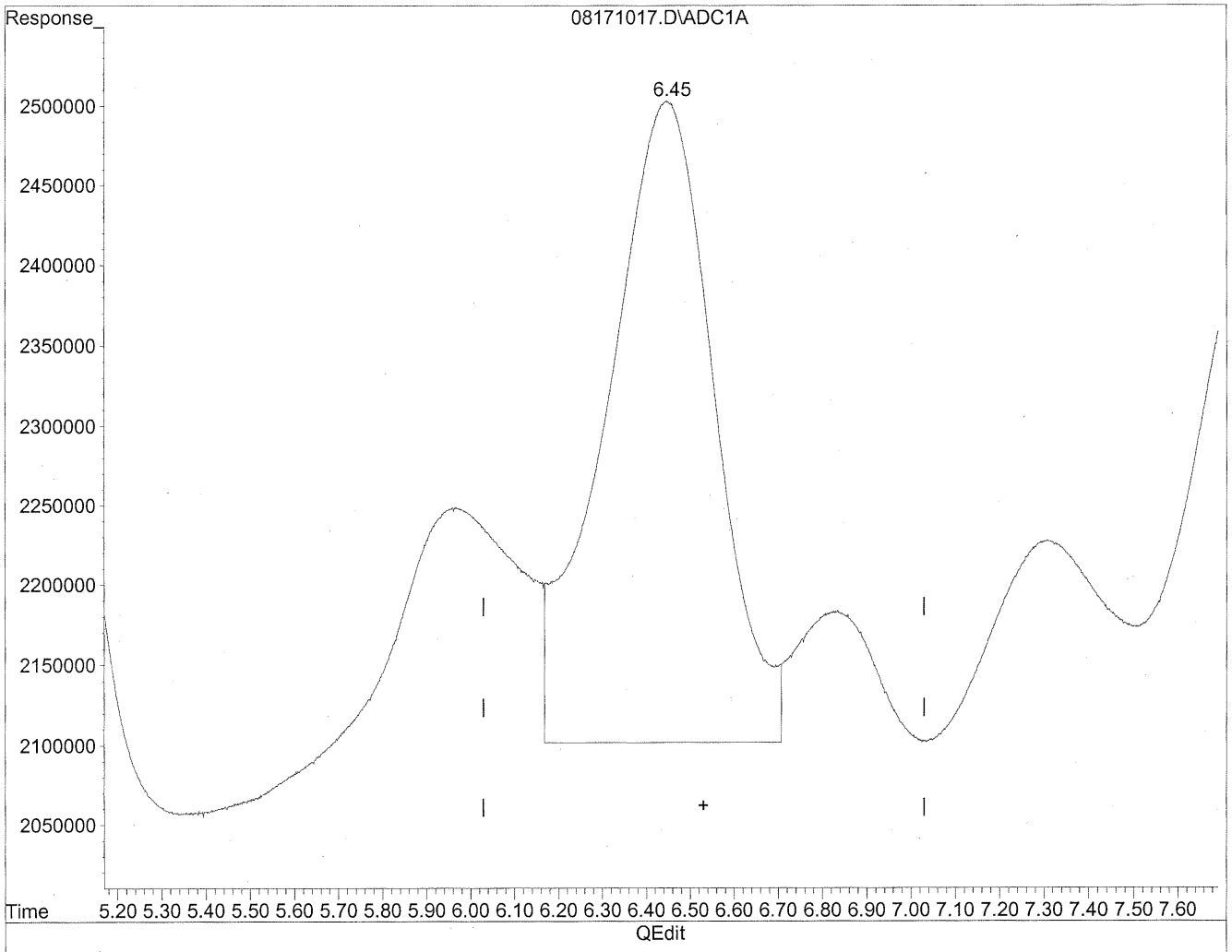
(+) = Expected Retention Time

08171017.D TO110709.M Sat Aug 22 14:31:45 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.45min 1038.895ng/ml m

response 68431305

HC 8/22/09
BC
KR 8/22/09

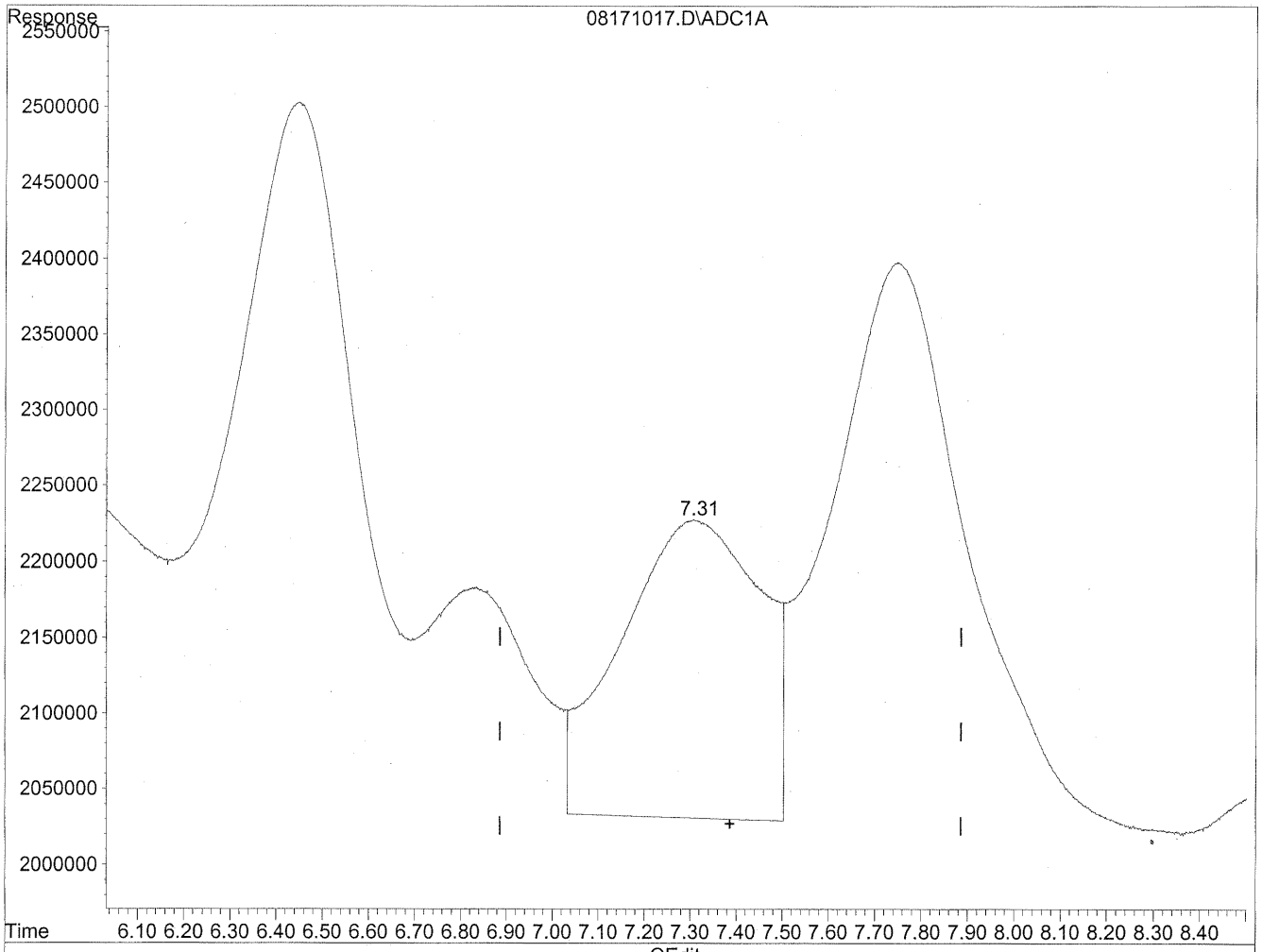
173

(+) = Expected Retention Time
08171017.D TO110709.M Sat Aug 22 14:32:04 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.31min 530.179ng/ml

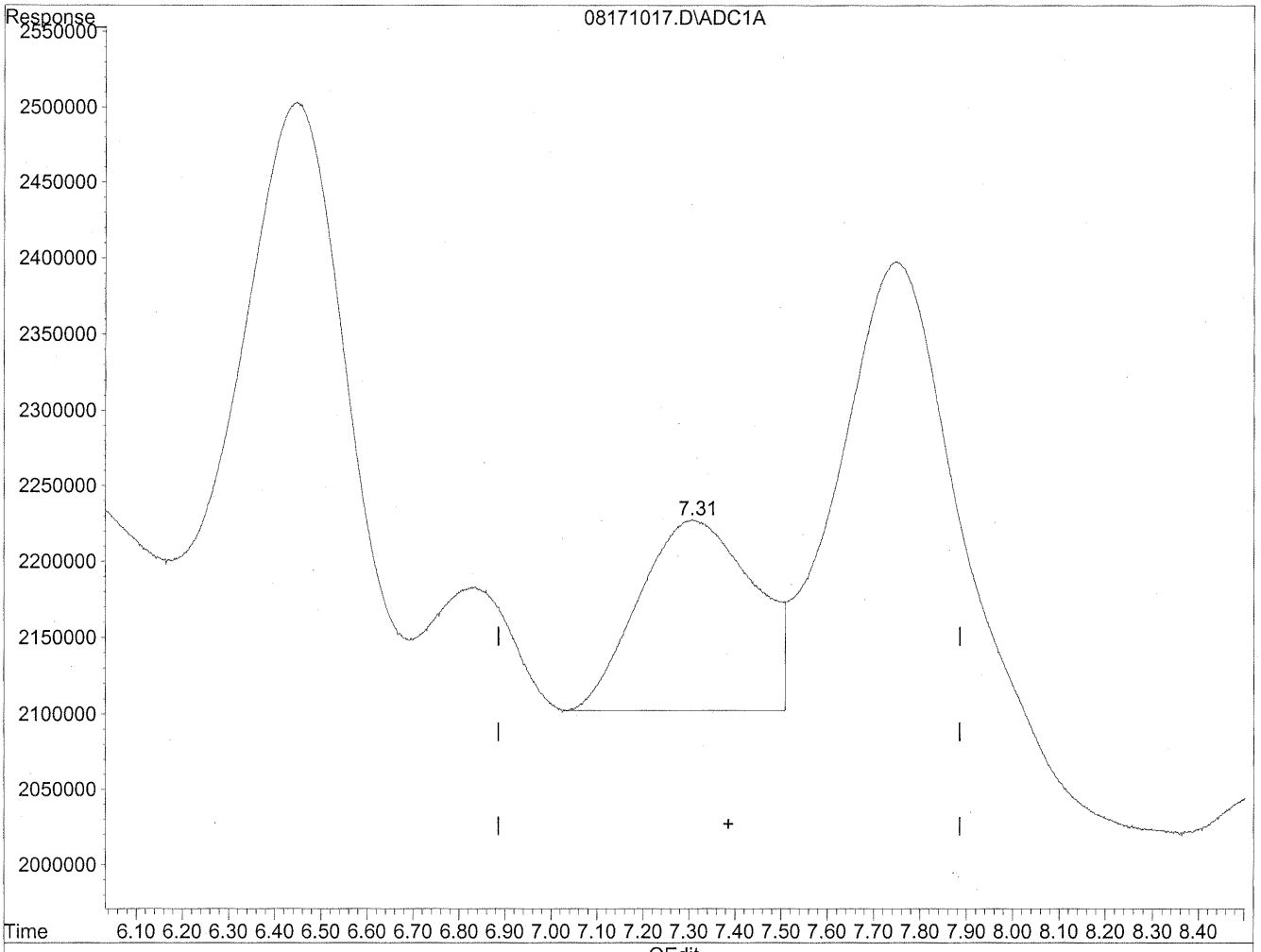
response 41487008

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.31min 277.520ng/ml m

response 21716251

HC 8/22/09

BC 8/23/09

175

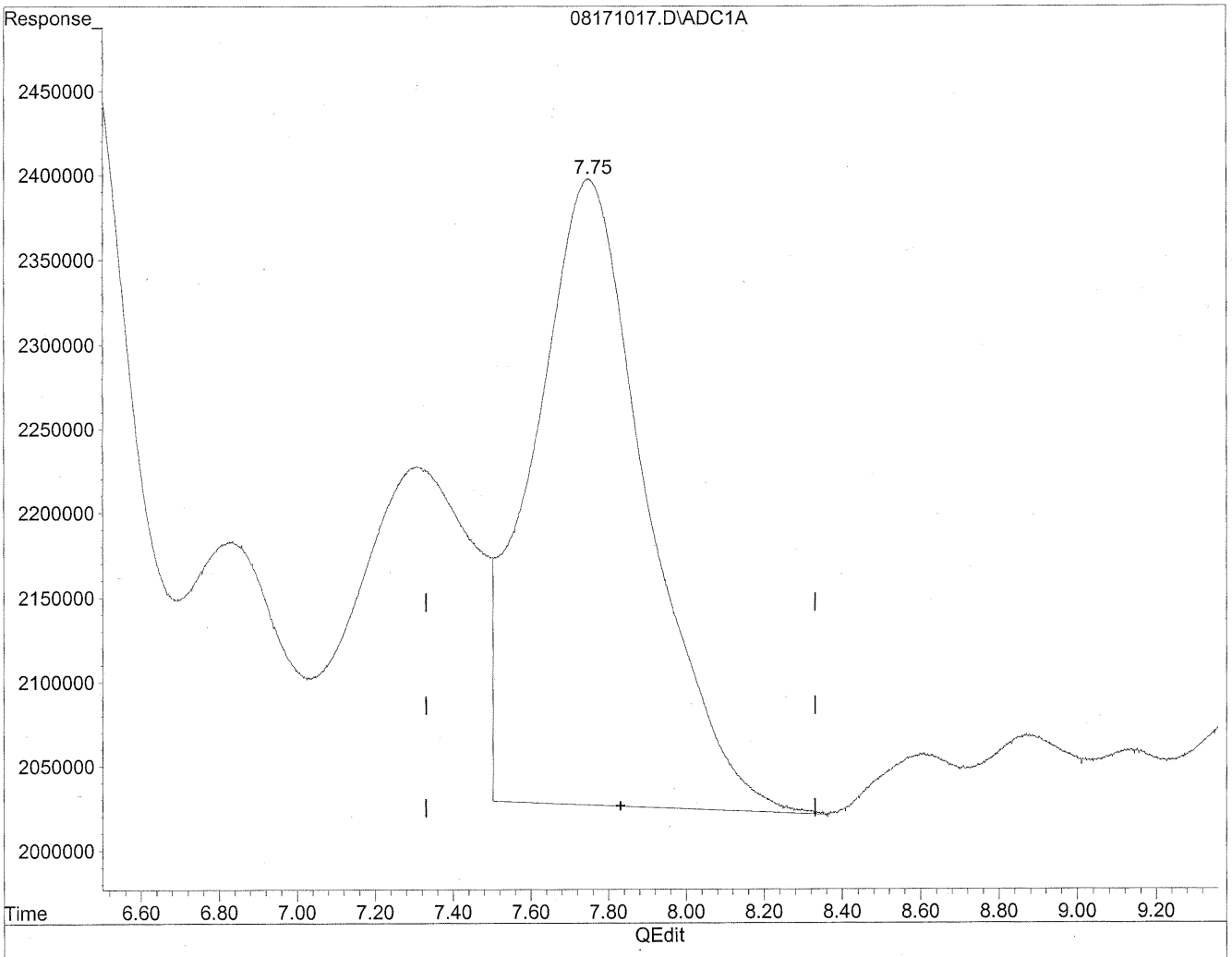
(+) = Expected Retention Time

08171017.D TO110709.M Sat Aug 22 14:32:17 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.75min 1030.486ng/ml

response 75745882

176

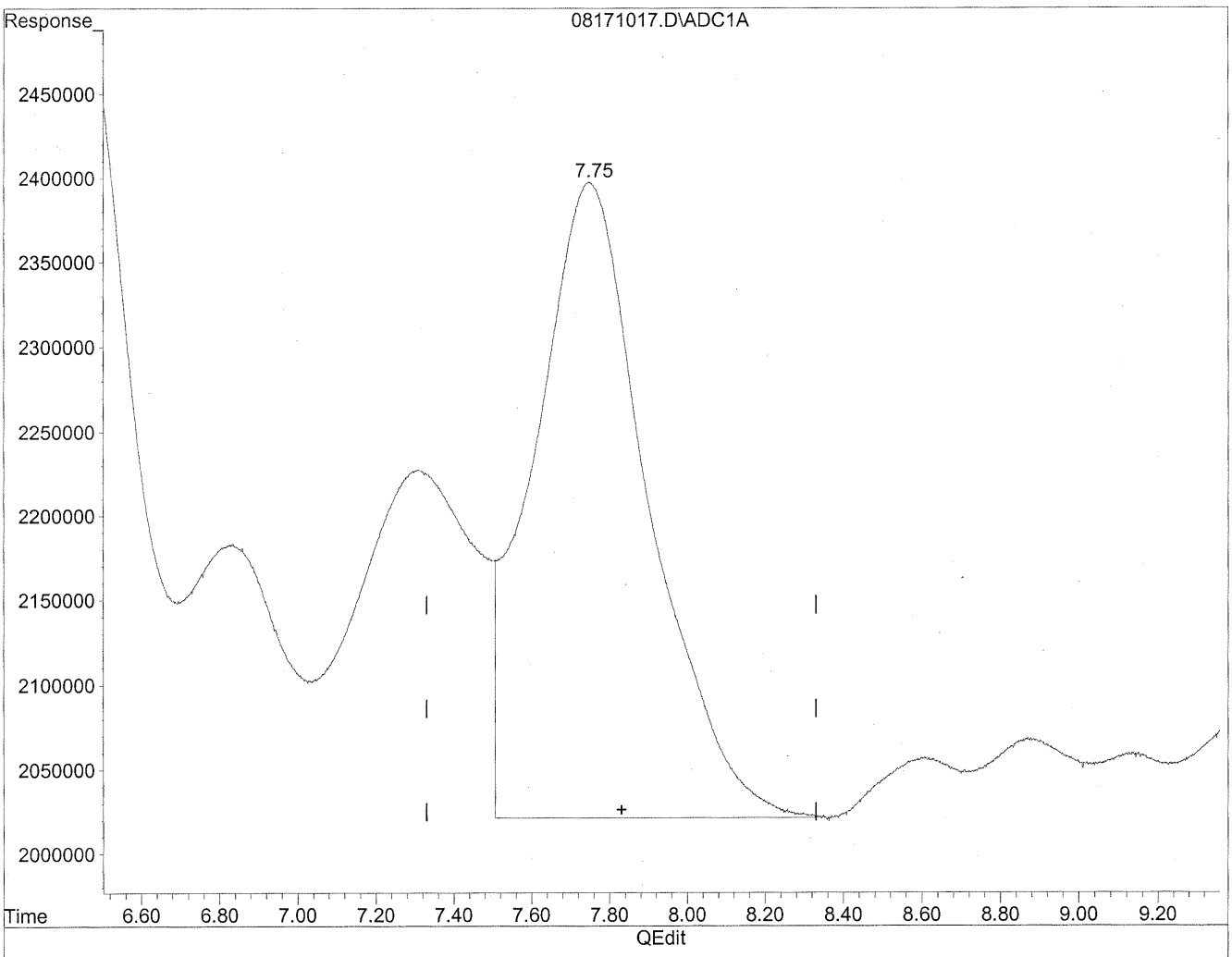
(+) = Expected Retention Time

08171017.D TO110709.M Sat Aug 22 14:32:22 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.75min 1050.376ng/ml m

response 77207887

HC 8/22/09
BC
KP 8/23/09

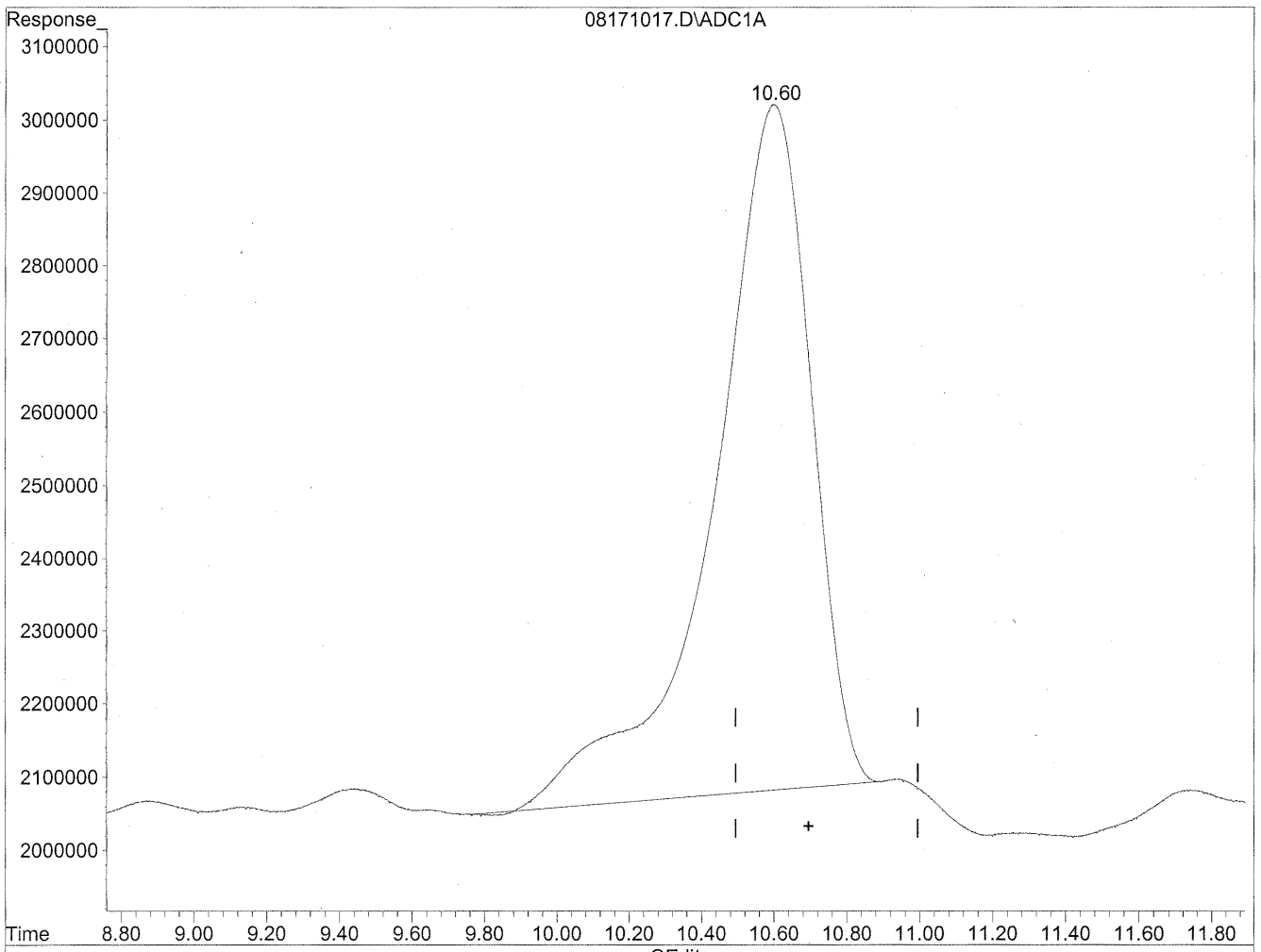
(+) = Expected Retention Time

08171017.D TO110709.M Sat Aug 22 14:32:33 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.60min 2601.018ng/ml

response 175162319

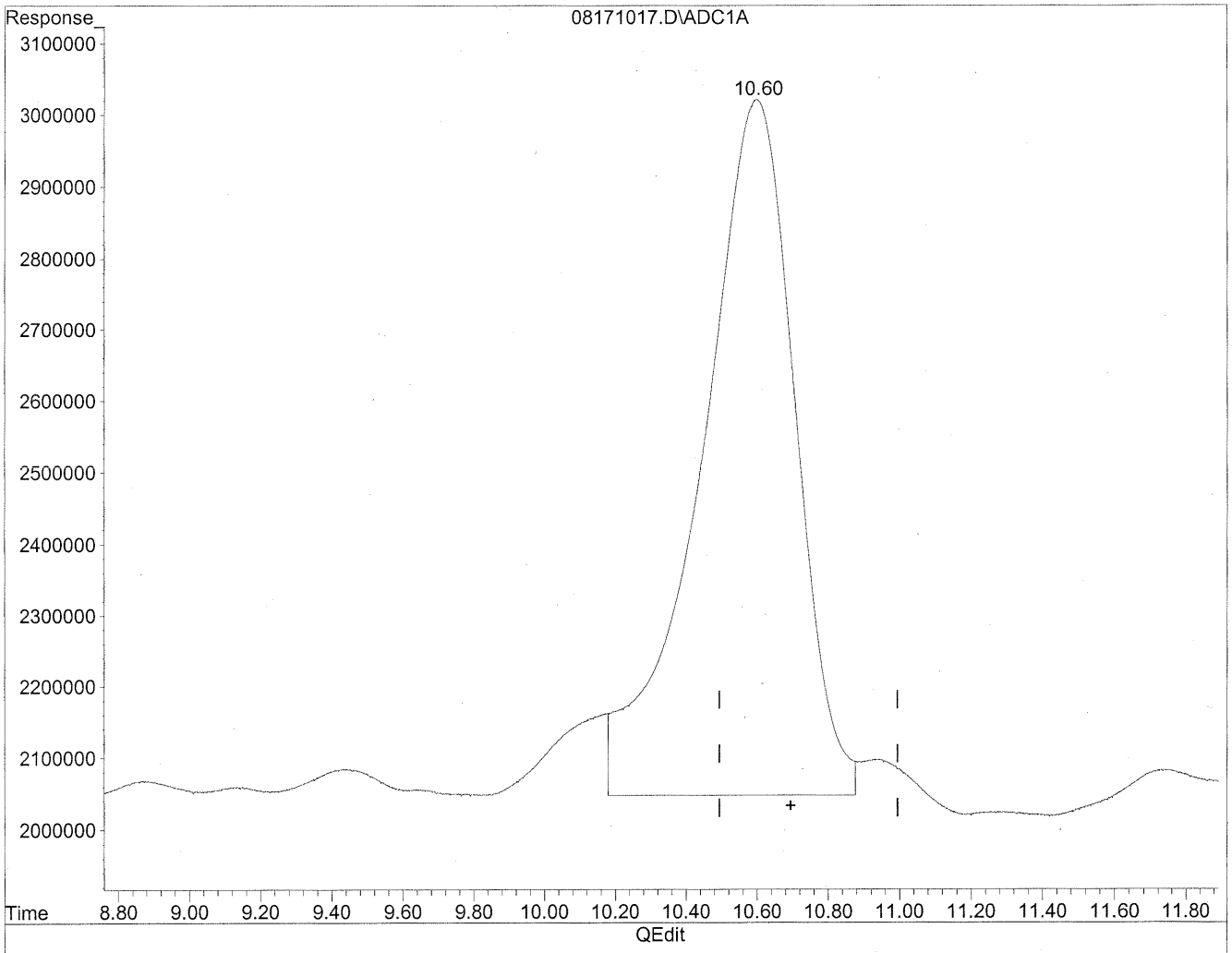
(+) = Expected Retention Time

08171017.D TO110709.M Sat Aug 22 14:32:38 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.60min 2655.258ng/ml m

response 178815052

HC 8/22/09 SH, BC

8/23/09

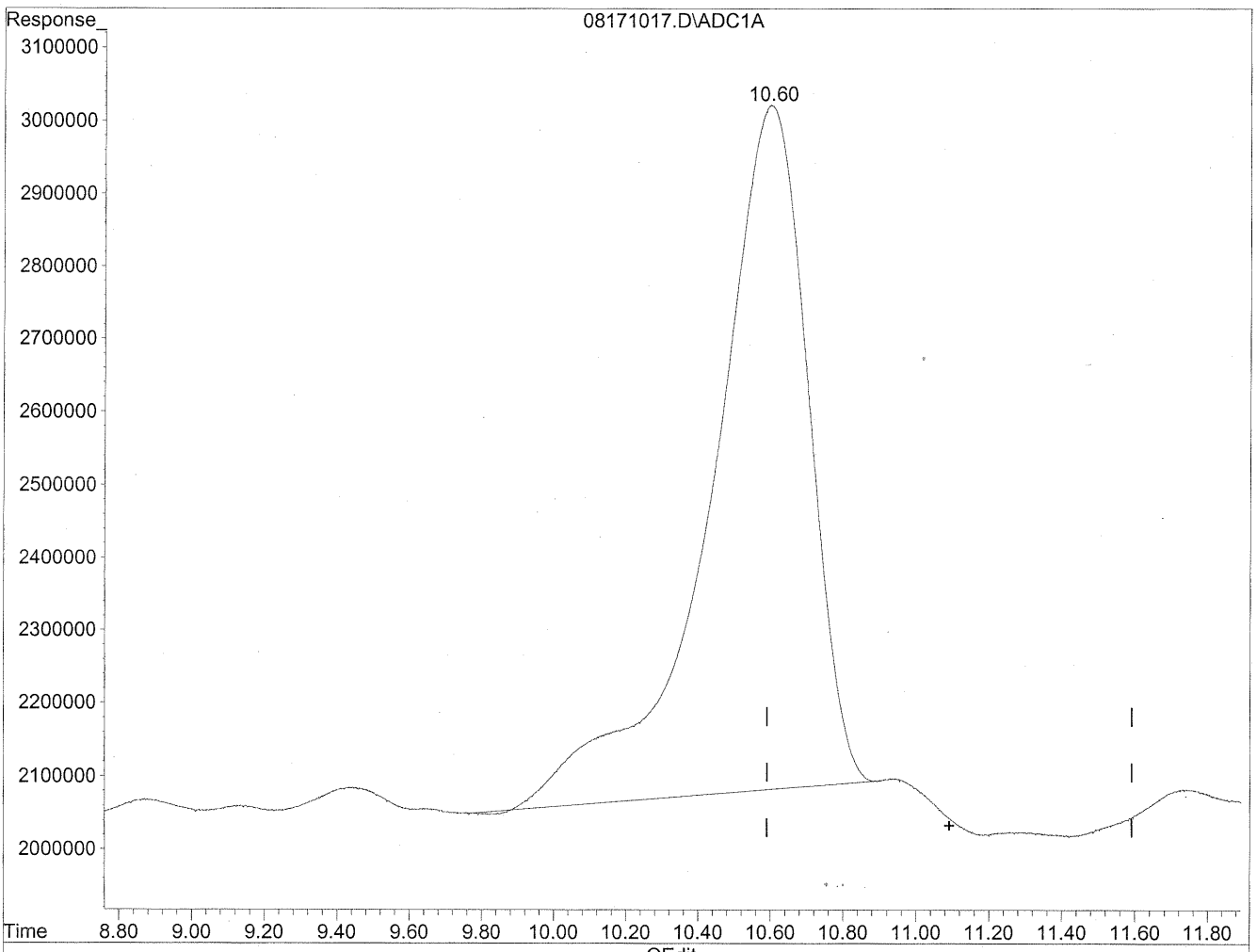
(+) = Expected Retention Time

08171017.D TO110709.M Sat Aug 22 14:32:52 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.60min 3573.765ng/ml

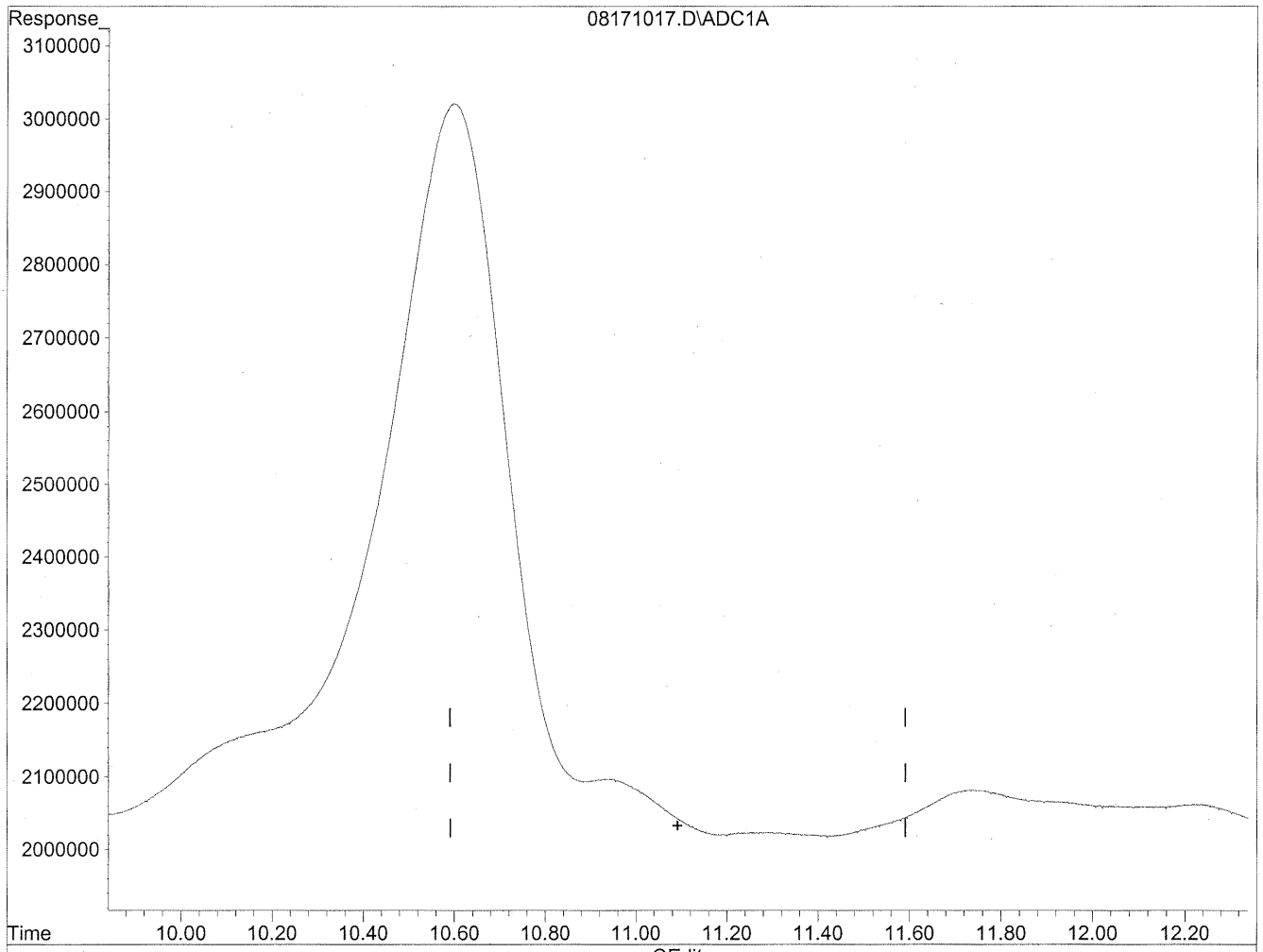
response 175162319

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171017.D Vial: 30
Acq On : 18 Aug 2009 7:54 pm Operator: HC
Sample : P0902772-007 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09
MP 8/23/09

181

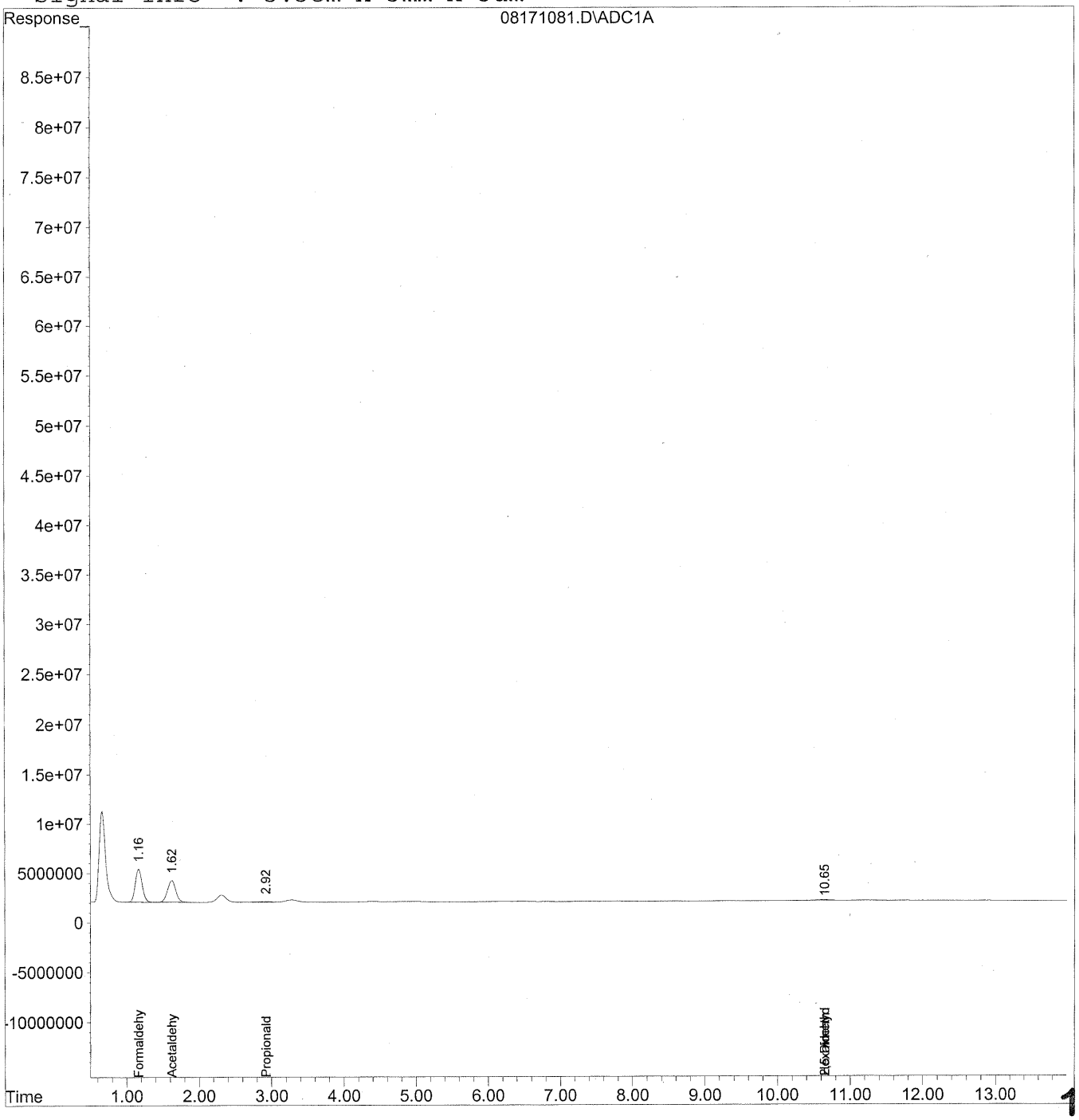
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171081.D Vial: 6
Acq On : 19 Aug 2009 11:57 am Operator: HC
Sample : P0902772-007 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:22 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



182

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171081.D Vial: 6
Acq On : 19 Aug 2009 11:57 am Operator: HC
Sample : P0902772-007 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:22 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

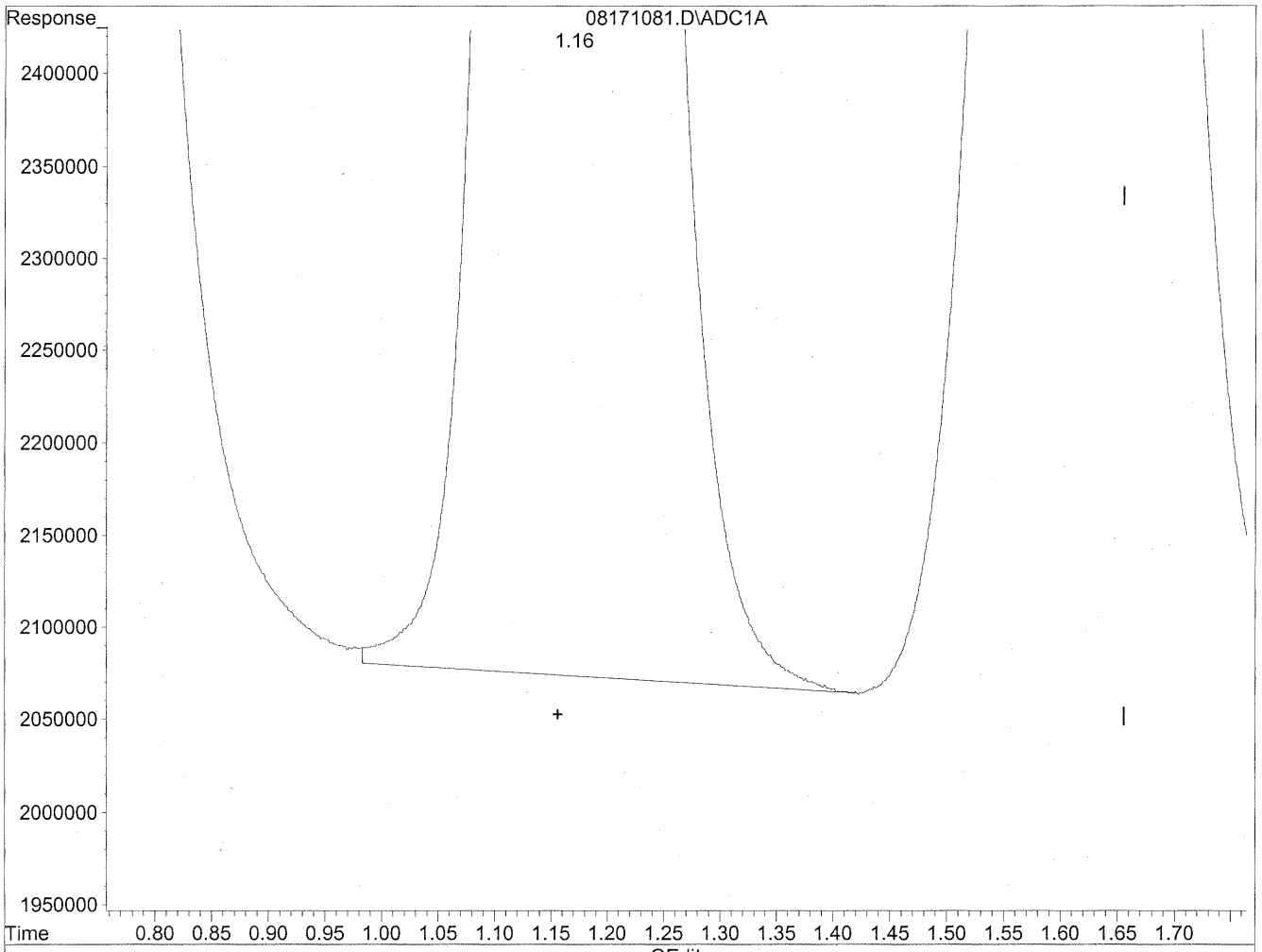
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	222271682	1210.753	ng/mlm
2)	Acetaldehyde	1.62	179724427	1281.700	ng/ml
3)	Propionaldehyde	2.92	9500502	89.043	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	10.64	15749175	233.862	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.64f	15749175	321.324	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171081.D Vial: 6
Acq On : 19 Aug 2009 11:57 am Operator: HC
Sample : P0902772-007 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 1217.232ng/ml

response 223461237

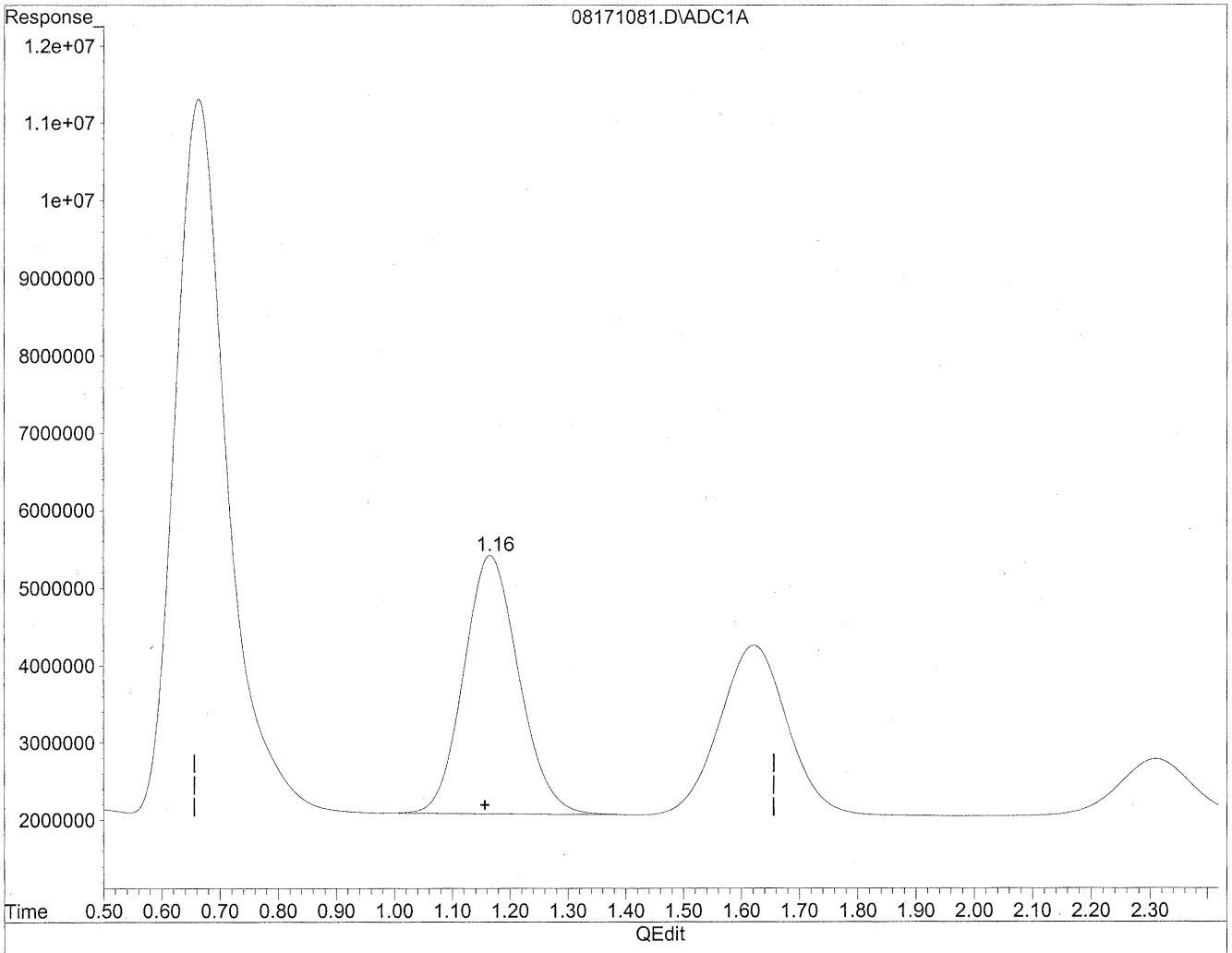
184

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171081.D Vial: 6
Acq On : 19 Aug 2009 11:57 am Operator: HC
Sample : P0902772-007 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.16min 1210.753ng/ml m

response 222271682

HC 8/22/09

K KP 8/23/09

185

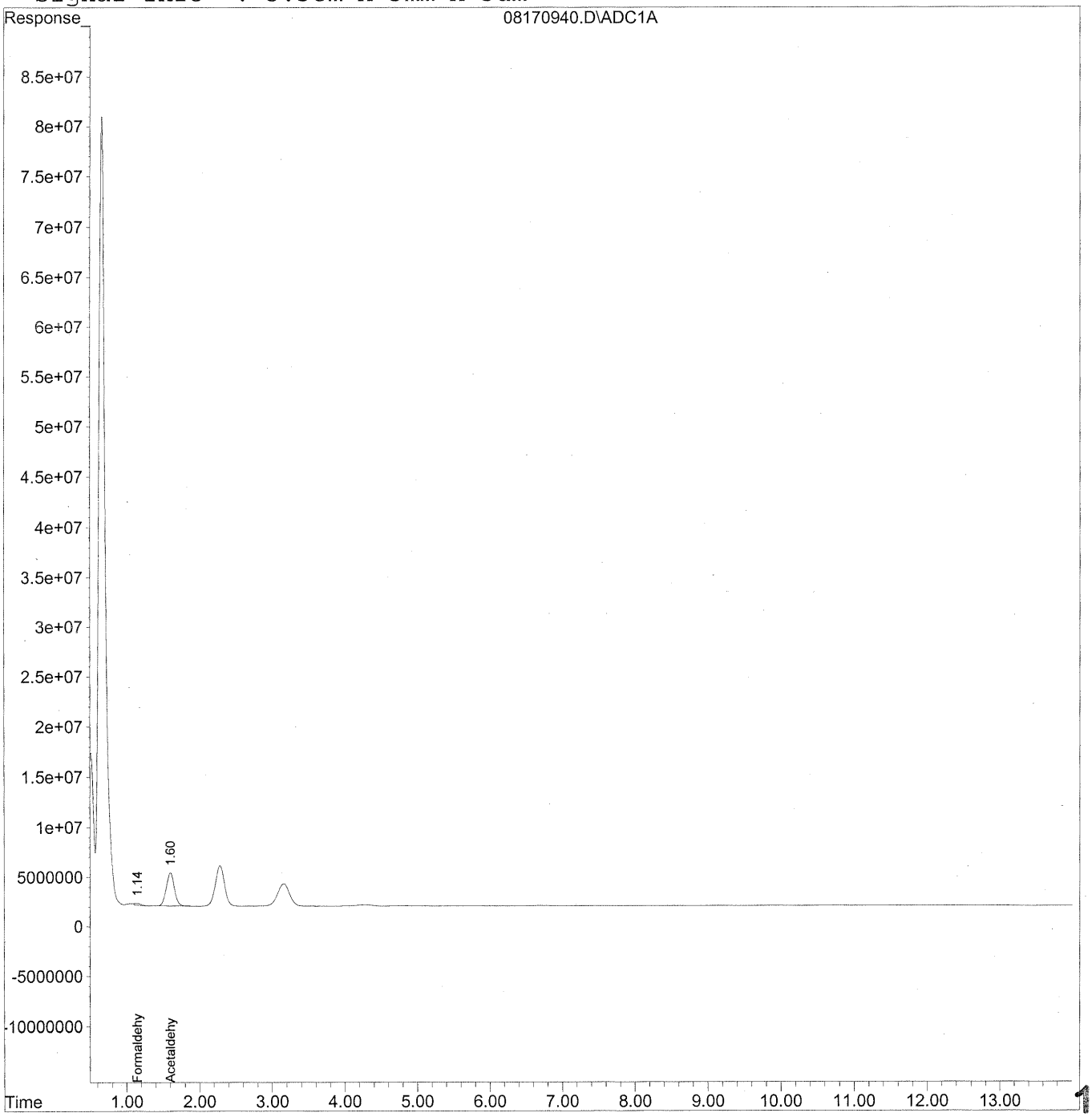
(+) = Expected Retention Time

08171081.D TO110709.M Sat Aug 22 17:22:16 2009

Data File : J:\LC01\DATA\TO11\2009_08\17\08170940.D Vial: 39
Acq On : 18 Aug 2009 12:36 am Operator: HC
Sample : P0902772-007 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:19 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



186

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170940.D Vial: 39
Acq On : 18 Aug 2009 12:36 am Operator: HC
Sample : P0902772-007 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:19 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

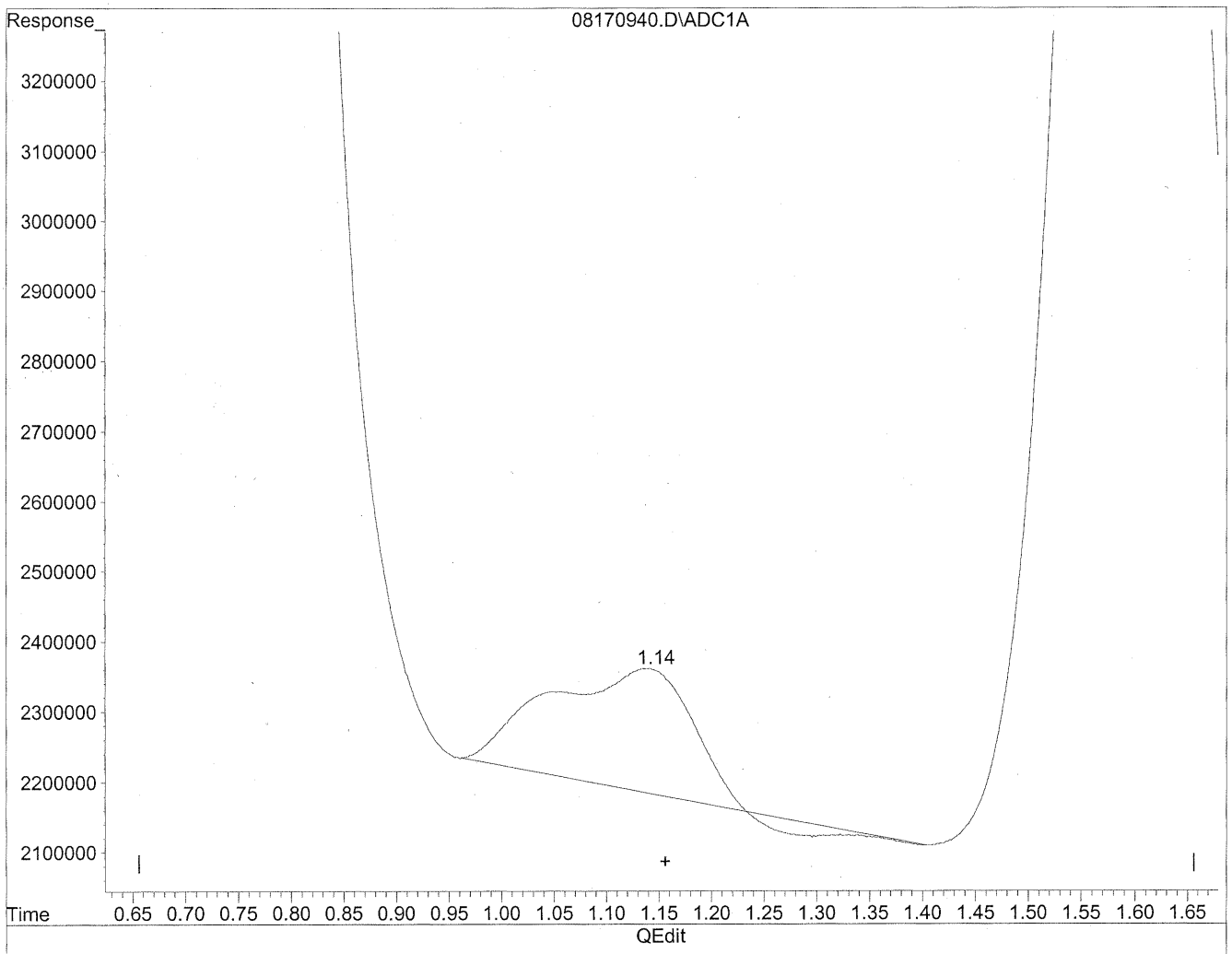
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.14	9897265	53.912	ng/mlm
2)	Acetaldehyde	1.60	261653338	1865.973	ng/ml
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170940.D Vial: 39
Acq On : 18 Aug 2009 12:36 am Operator: HC
Sample : P0902772-007 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:18 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.14min 83.339ng/ml

response 15299466

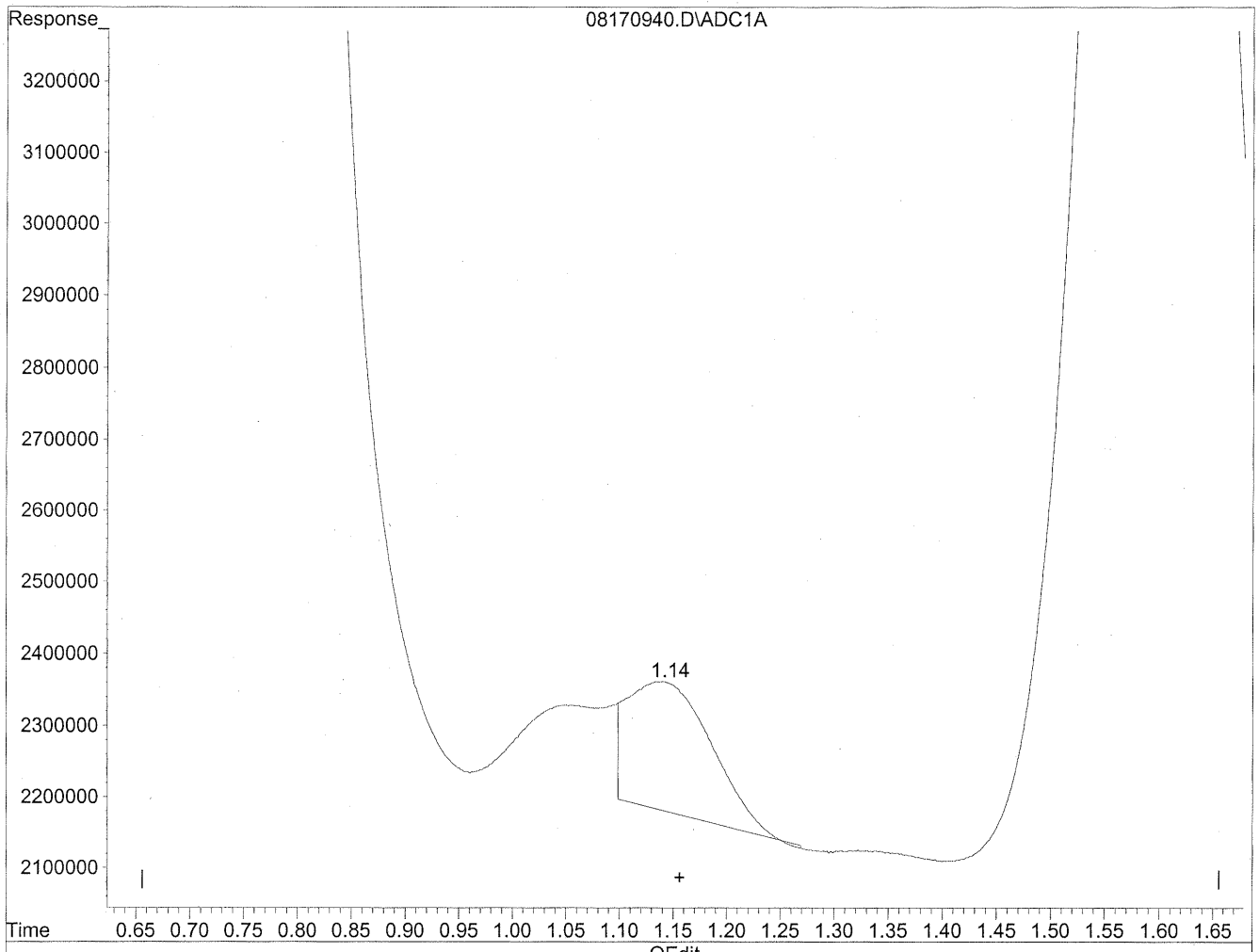
188

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170940.D Vial: 39
Acq On : 18 Aug 2009 12:36 am Operator: HC
Sample : P0902772-007 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:18 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.14min 53.912ng/ml m

response 9897265

HC 8/21/09 SP

KF 8/22/09

189

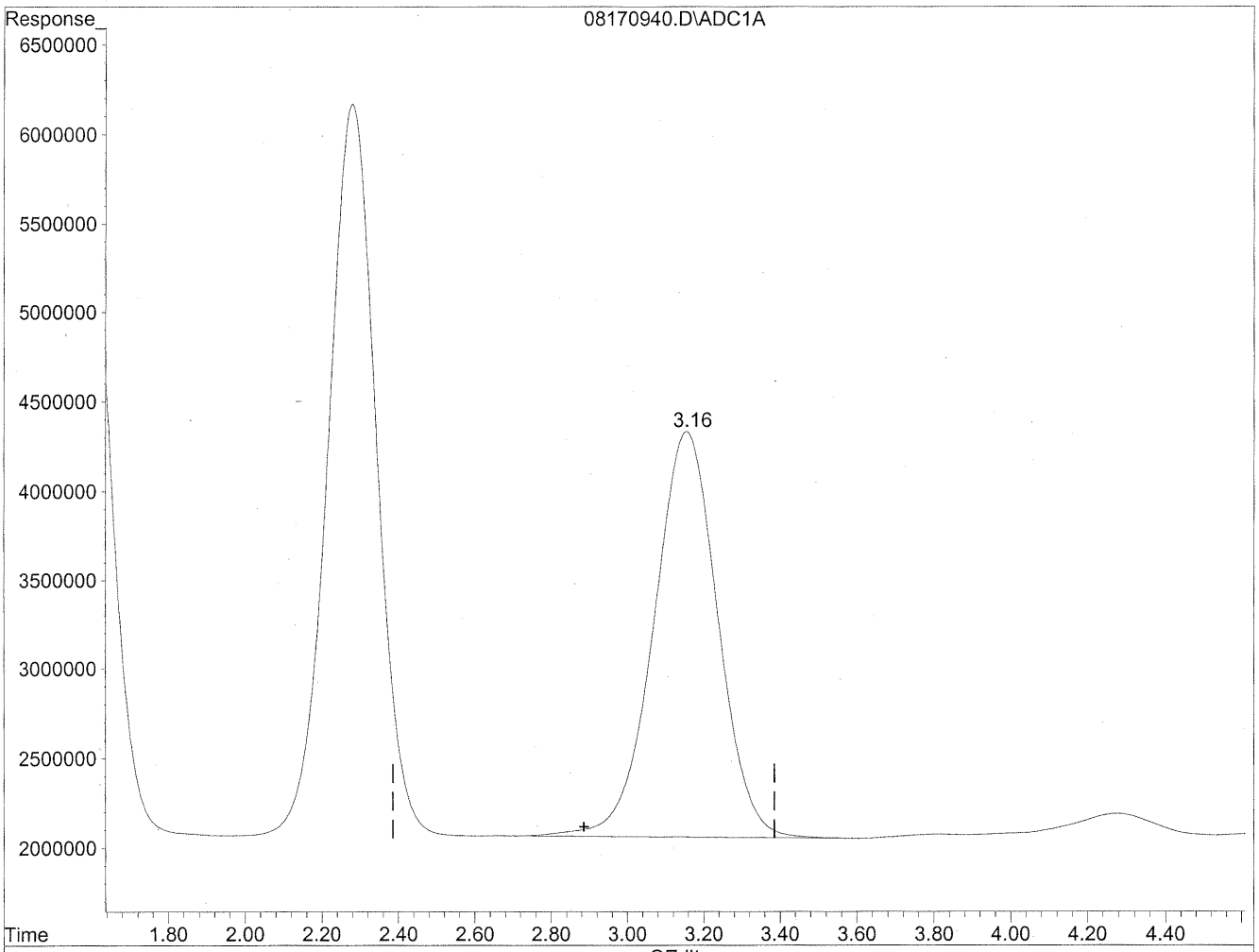
(+) = Expected Retention Time

08170940.D TO110709.M Fri Aug 21 18:18:58 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170940.D Vial: 39
Acq On : 18 Aug 2009 12:36 am Operator: HC
Sample : P0902772-007 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:18 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.15min 2484.258ng/ml

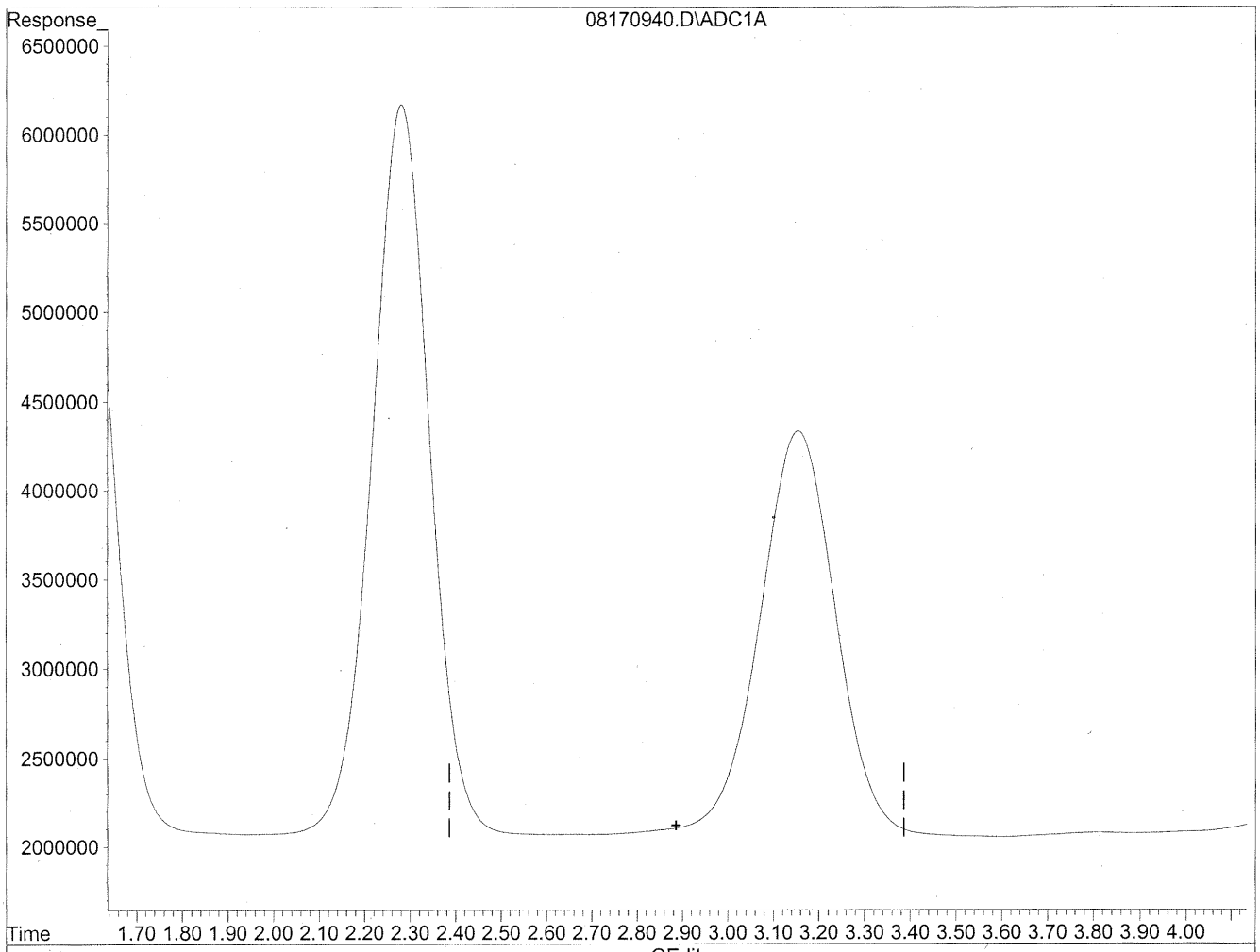
response 265058483

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170940.D Vial: 39
Acq On : 18 Aug 2009 12:36 am Operator: HC
Sample : P0902772-007 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:18 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/21/09 MP

KH 8/23/09

191

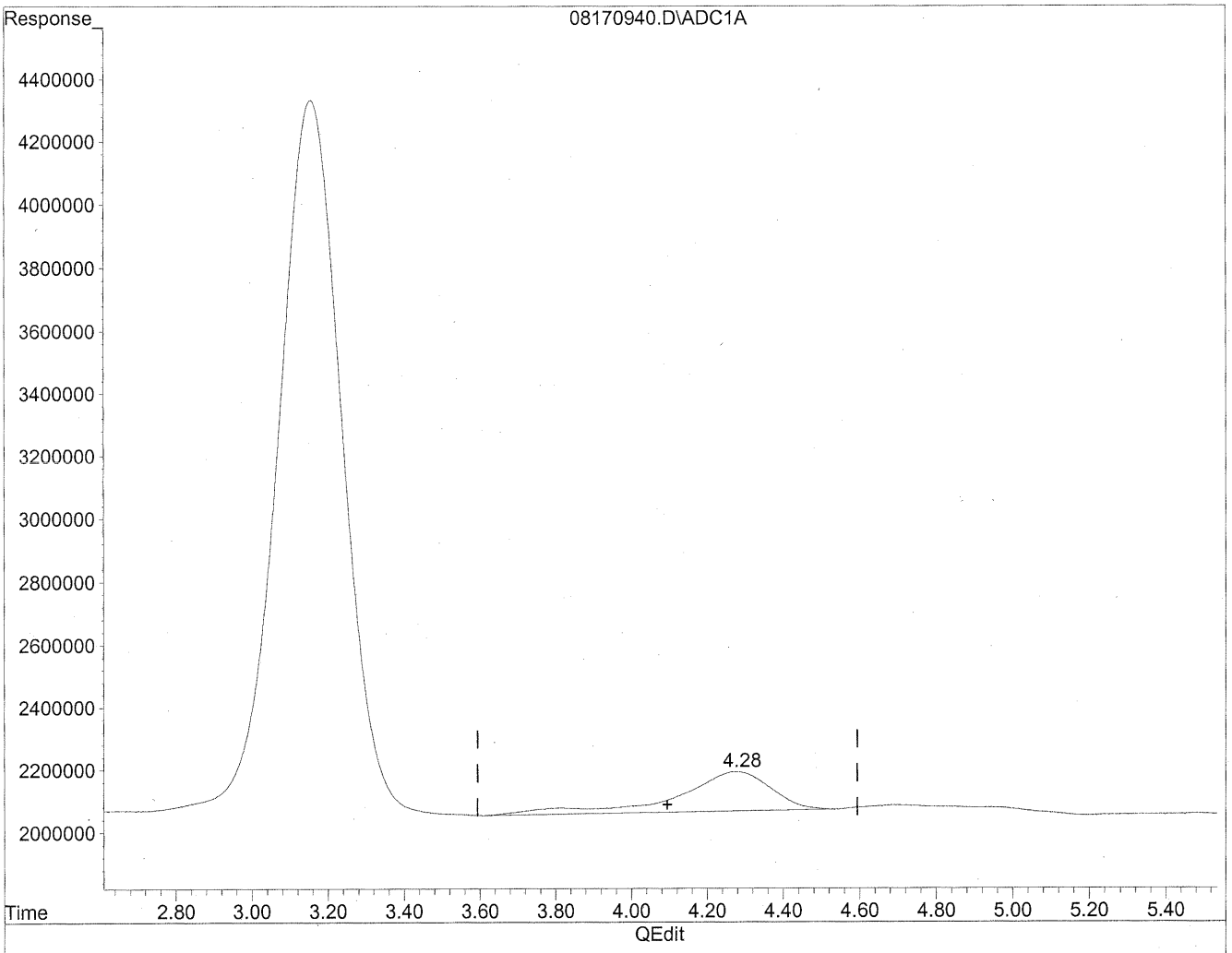
(+) = Expected Retention Time

08170940.D TO110709.M Fri Aug 21 18:19:09 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170940.D Vial: 39
Acq On : 18 Aug 2009 12:36 am Operator: HC
Sample : P0902772-007 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:18 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.28min 223.572ng/ml

response 21779271

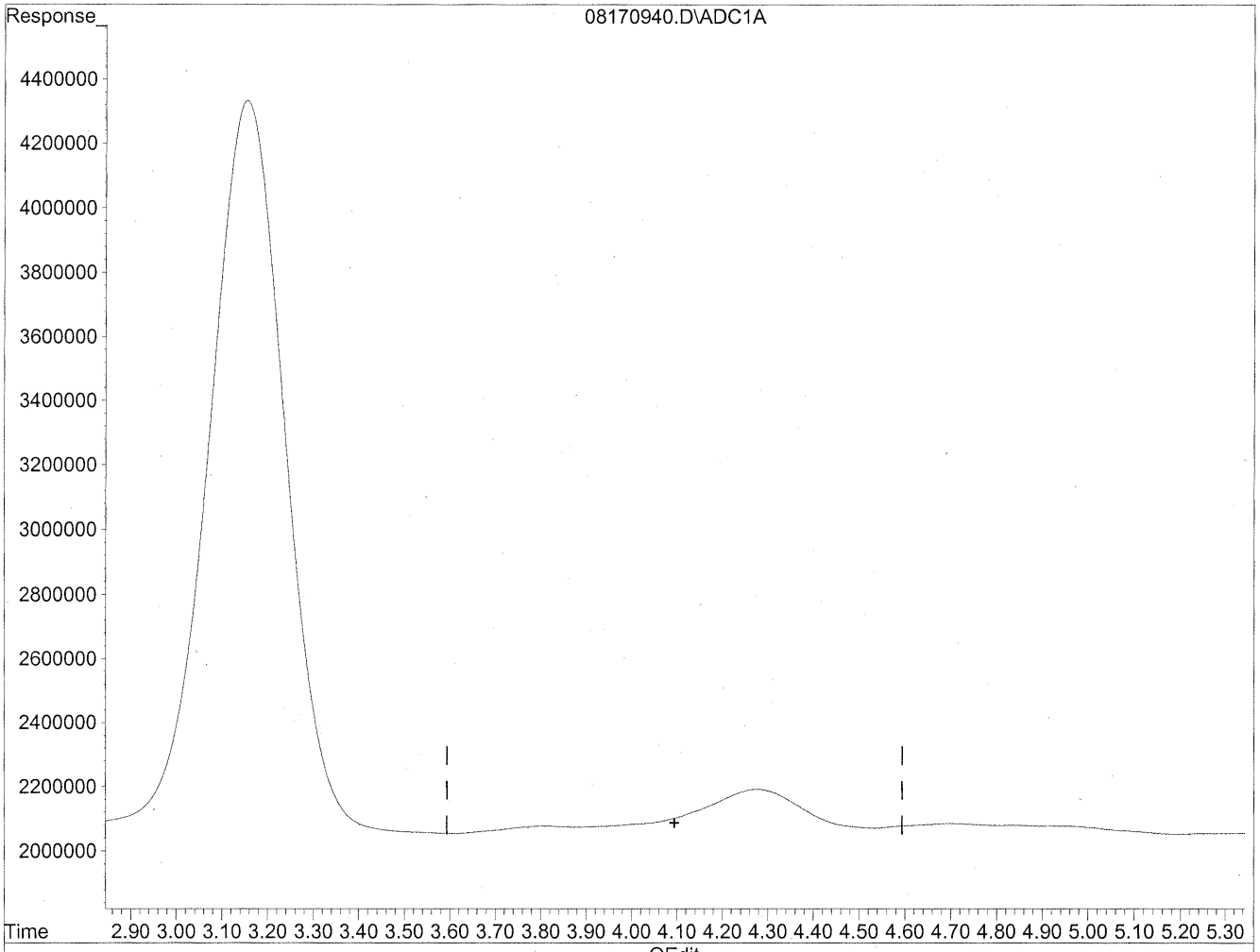
(+) = Expected Retention Time

08170940.D TO110709.M Fri Aug 21 18:19:13 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170940.D Vial: 39
Acq On : 18 Aug 2009 12:36 am Operator: HC
Sample : P0902772-007 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:18 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

hc 8/21/09 mp

hp 8/23/09

193

(+) = Expected Retention Time
08170940.D TO110709.M Fri Aug 21 18:19:16 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **100444**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P0902772-008

Test Code: EPA Method TO-11A
Instrument ID: Waters LC Module I Plus/UV_Vis 360/LC1
Analyst: Hani Cherazaie
Sampling Media: Silica Gel DNPH Tube
Test Notes: **BC**

Date Collected: 8/10/09
Date Received: 8/12/09
Date Analyzed: 8/18 - 8/19/09
Desorption Volume: 1.0 ml
Volume Sampled: 108.2 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**12,000**	**110**	0.92	**91**	0.75	
75-07-0	**Acetaldehyde**	**15,000**	**140**	0.92	**79**	0.51	**BT**
123-38-6	**Propionaldehyde**	**880**	**8.1**	0.92	**3.4**	0.39	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.92	ND	0.32	
123-72-8	**Butyraldehyde**	**540**	**5.0**	0.92	**1.7**	0.31	
100-52-7	**Benzaldehyde**	**1,200**	**11**	0.92	**2.5**	0.21	
590-86-3	**Isovaleraldehyde**	**390**	**3.6**	0.92	**1.0**	0.26	
110-62-3	**Valeraldehyde**	**1,400**	**13**	0.92	**3.8**	0.26	
529-20-4	o-Tolualdehyde	< 100	ND	0.92	ND	0.19	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	1.8	ND	0.38	
66-25-1	**n-Hexaldehyde**	**3,100**	**28**	0.92	**7.0**	0.23	
5779-94-2	**2,5-Dimethylbenzaldehyde**	**230**	**2.1**	0.92	**0.39**	0.17	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
BT = Results indicated possible breakthrough; back section > 10% front section.

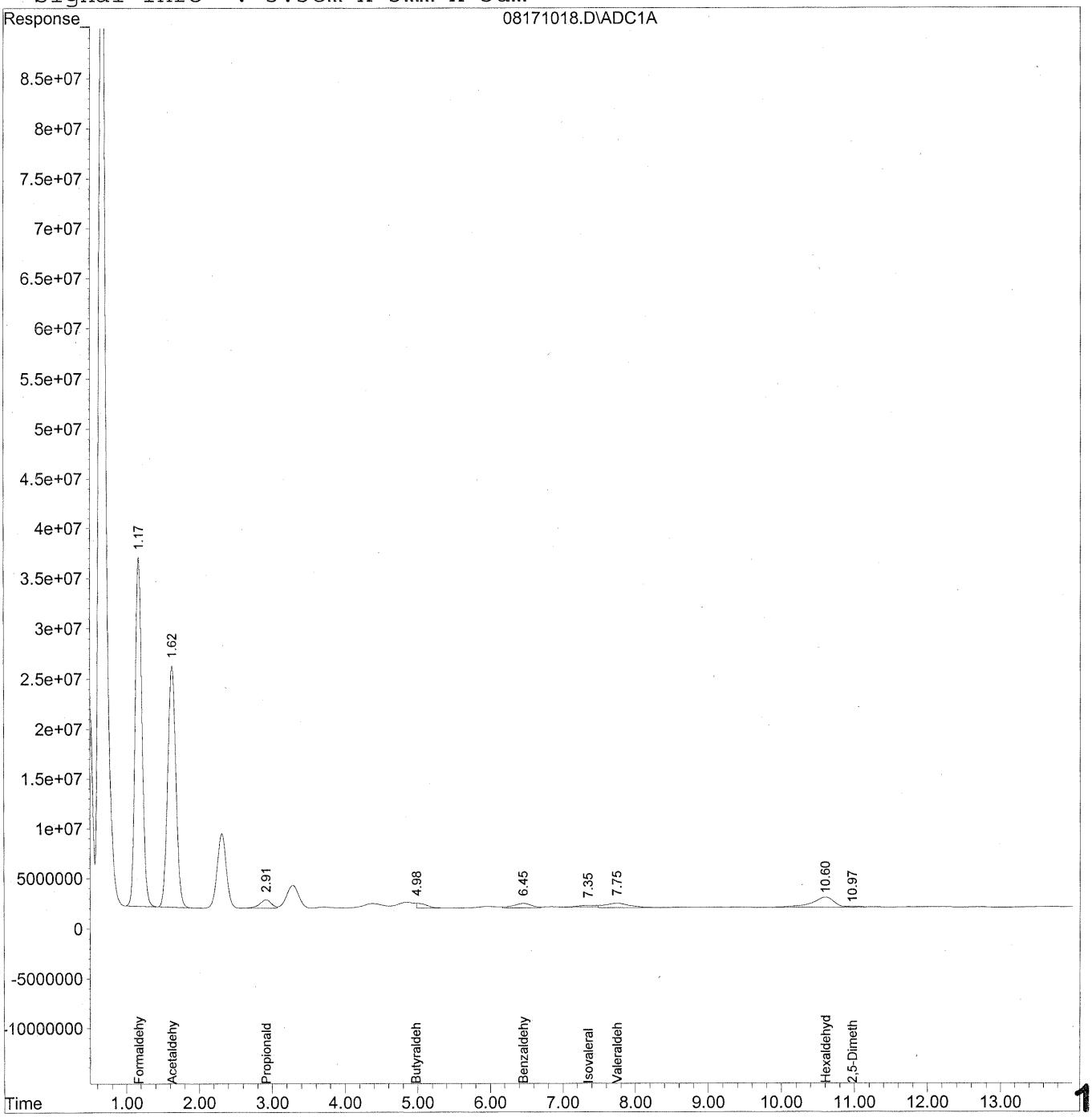
Verified By: _____ Date: 8/25/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171018.D Vial: 31
Acq On : 18 Aug 2009 8:09 pm Operator: HC
Sample : P0902772-008 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:35 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171018.D Vial: 31
Acq On : 18 Aug 2009 8:09 pm Operator: HC
Sample : P0902772-008 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:35 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

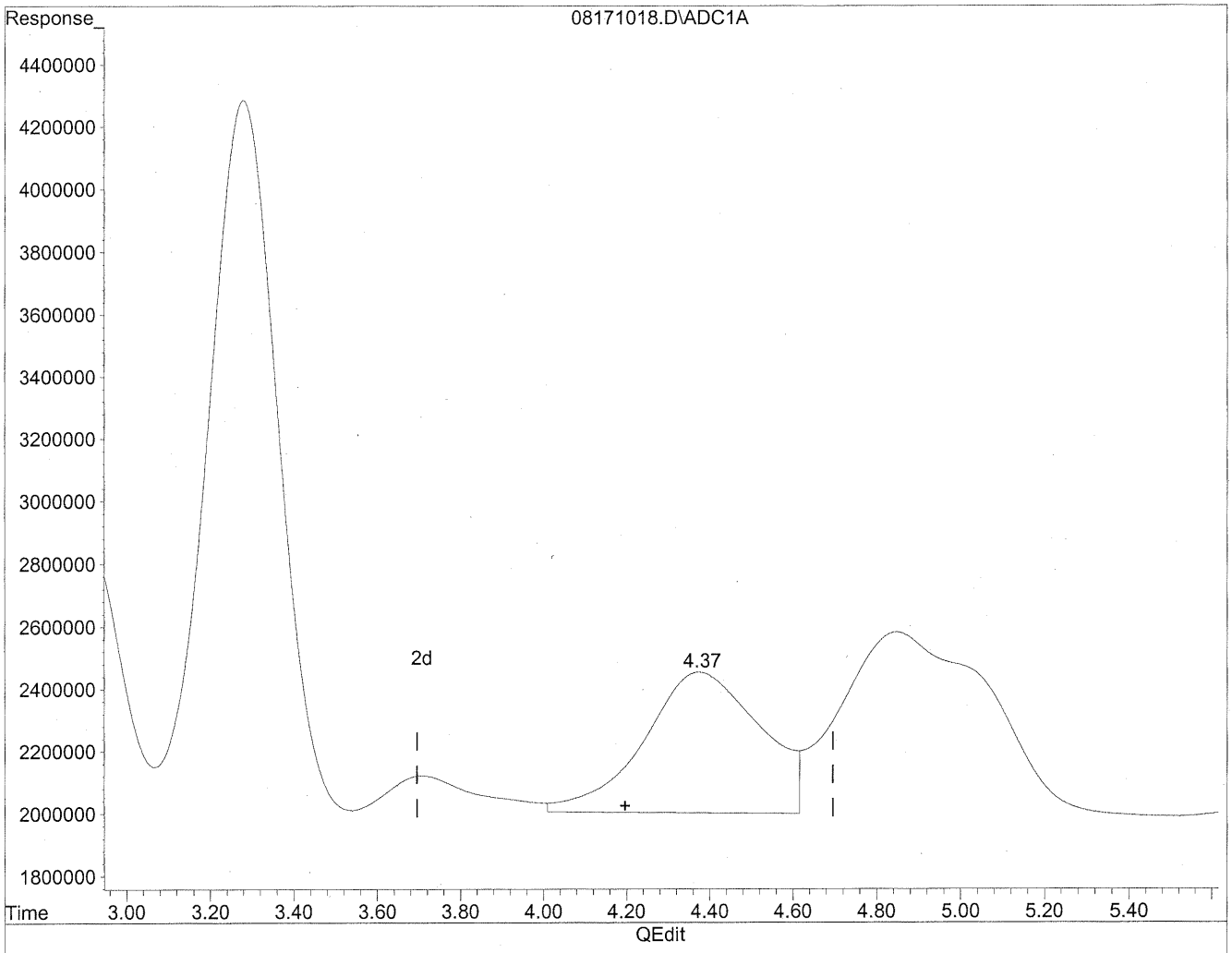
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	2277507994	12405.985	ng/ml
2)	Acetaldehyde	1.62	1917989658	13678.089	ng/ml
3)	Propionaldehyde	2.91	94033909	881.332	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	4.98	47466121	537.335	ng/mlm
6)	Benzaldehyde	6.46	77908191	1182.769	ng/ml
7)	Isovaleraldehyde	7.35	30782051	393.376	ng/mlm
8)	Valeraldehyde	7.75	106456531	1448.289	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	10.60	207572245	3082.279	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.97	11300857	230.567	ng/mlm

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171018.D Vial: 31
Acq On : 18 Aug 2009 8:09 pm Operator: HC
Sample : P0902772-008 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.37min 904.193ng/ml

response 88082158

197

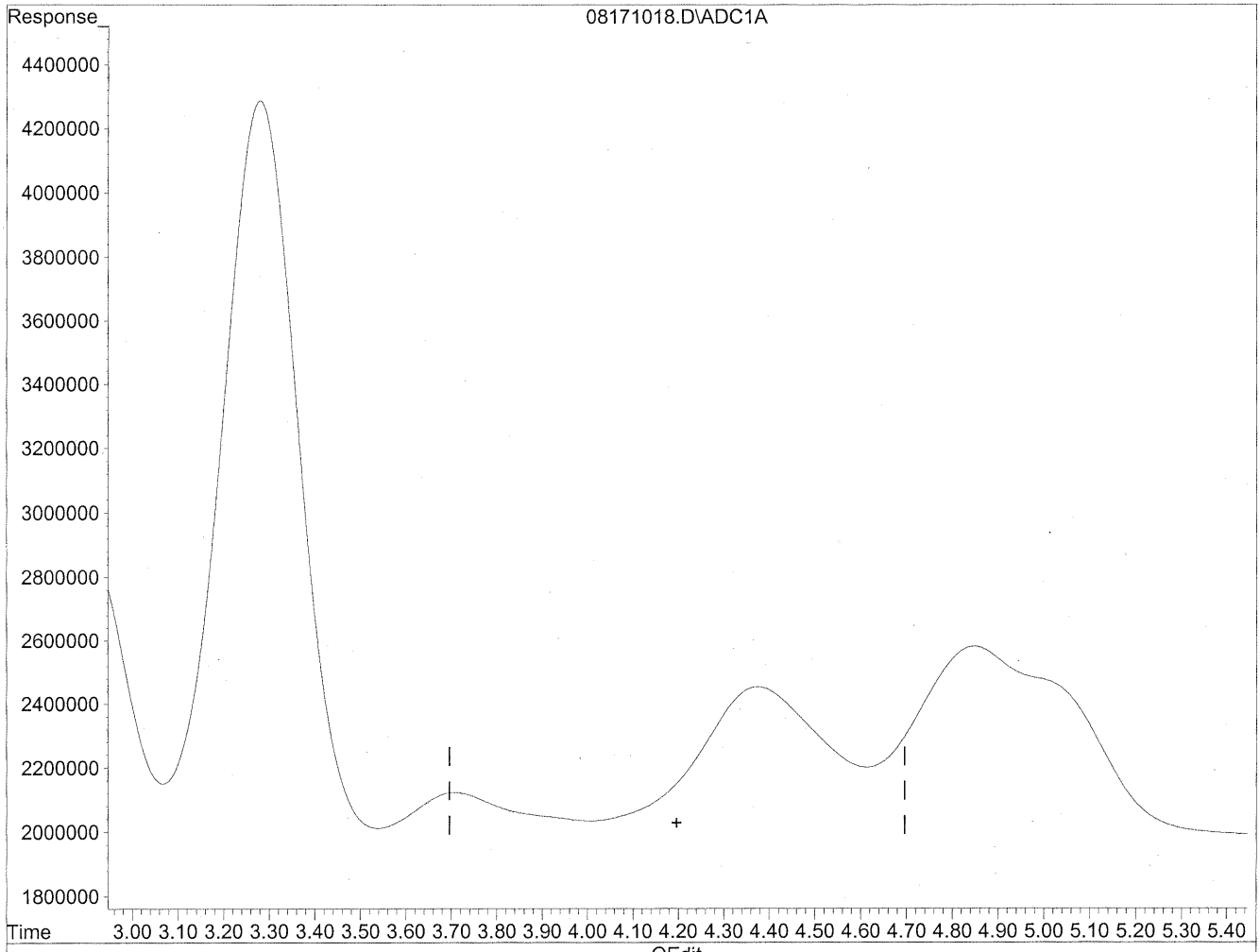
(+) = Expected Retention Time

08171018.D TO110709.M Sat Aug 22 14:33:42 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171018.D Vial: 31
Acq On : 18 Aug 2009 8:09 pm Operator: HC
Sample : P0902772-008 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HL 8/22/09 MP

8/23/09

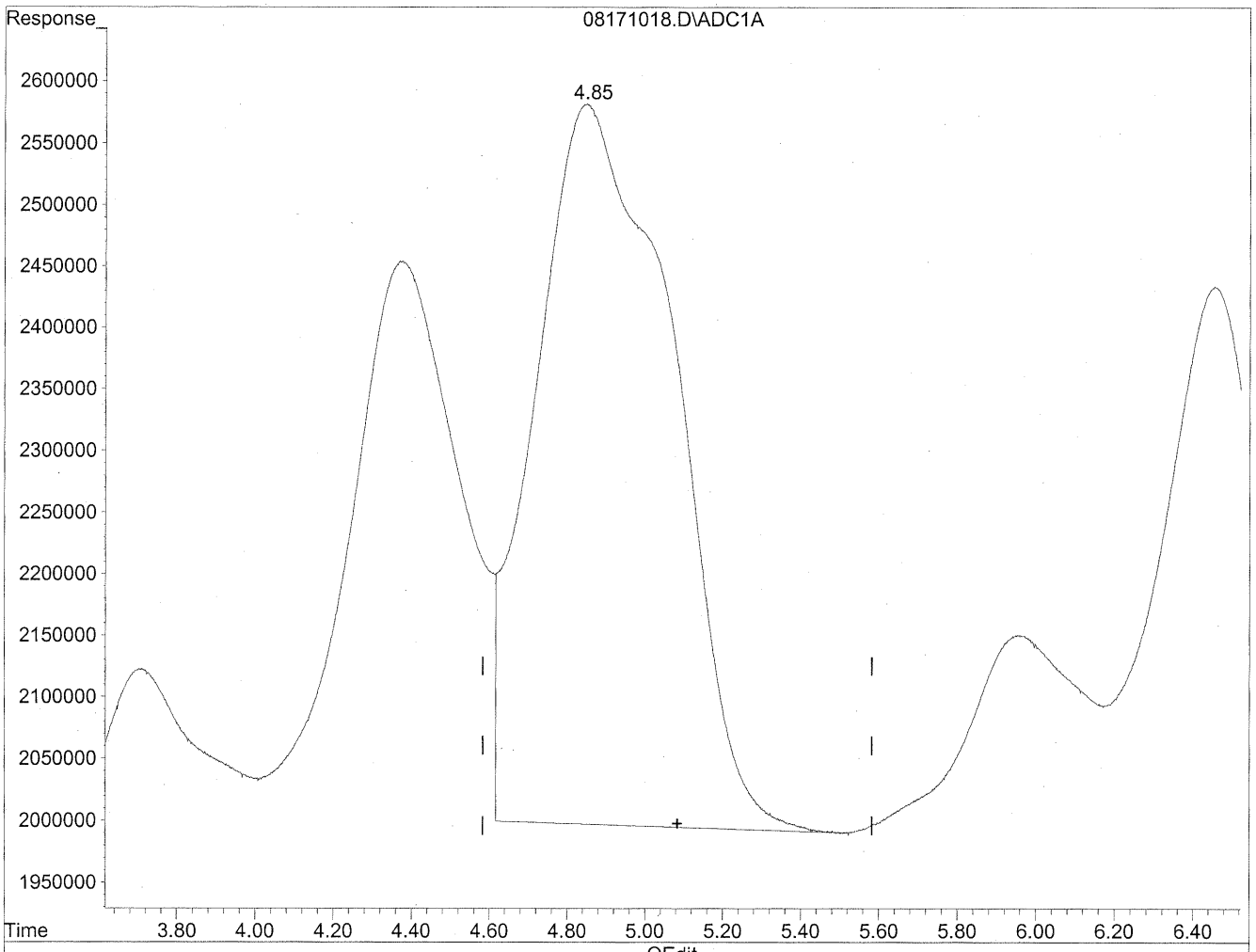
198

(+) = Expected Retention Time
08171018.D TO110709.M Sat Aug 22 14:33:45 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171018.D Vial: 31
Acq On : 18 Aug 2009 8:09 pm Operator: HC
Sample : P0902772-008 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.85min 1630.216ng/ml

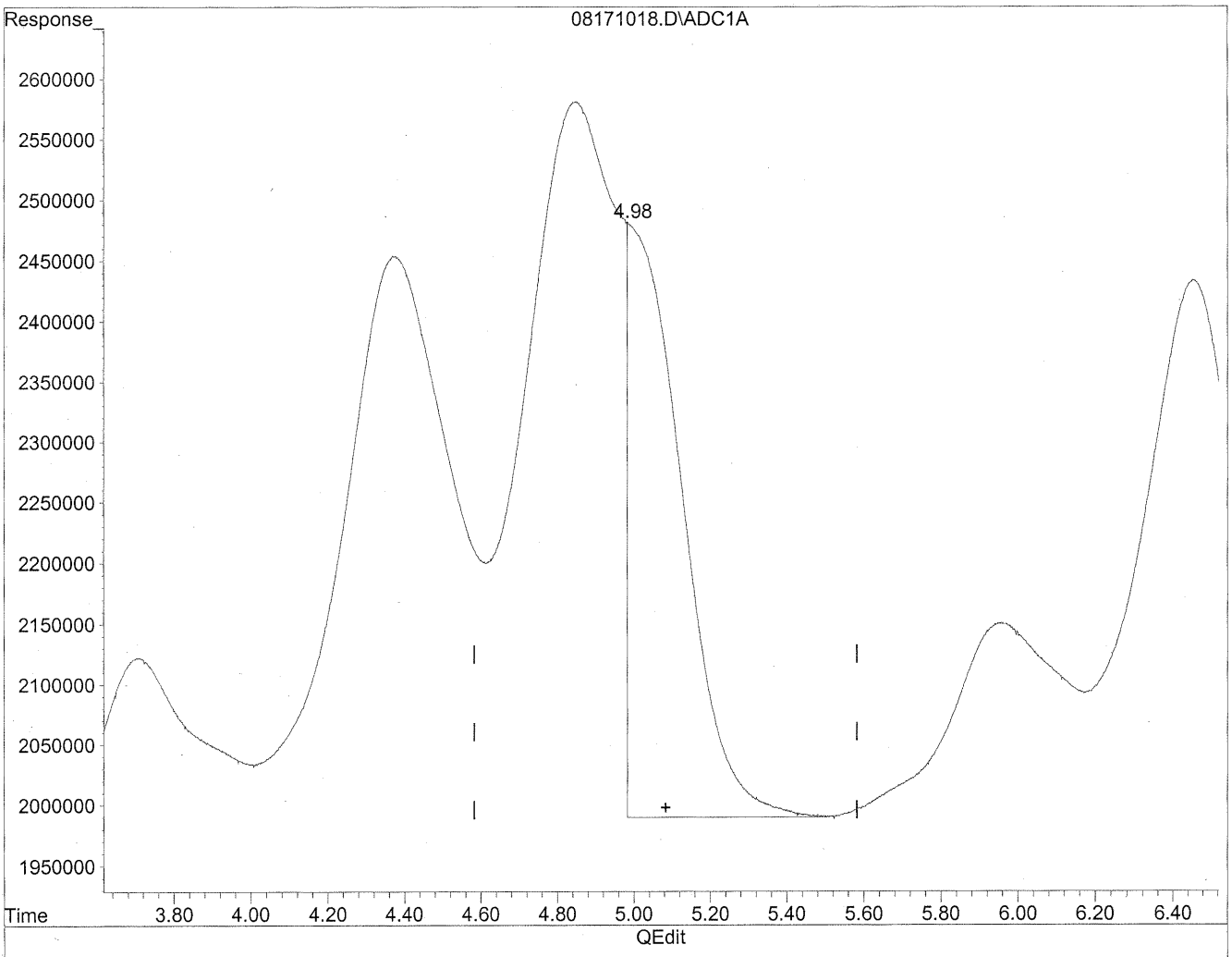
response 144007004

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171018.D Vial: 31
Acq On : 18 Aug 2009 8:09 pm Operator: HC
Sample : P0902772-008 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.98min 537.335ng/ml m

response 47466121

hc 8/22/09
SP
KP 8/23/09

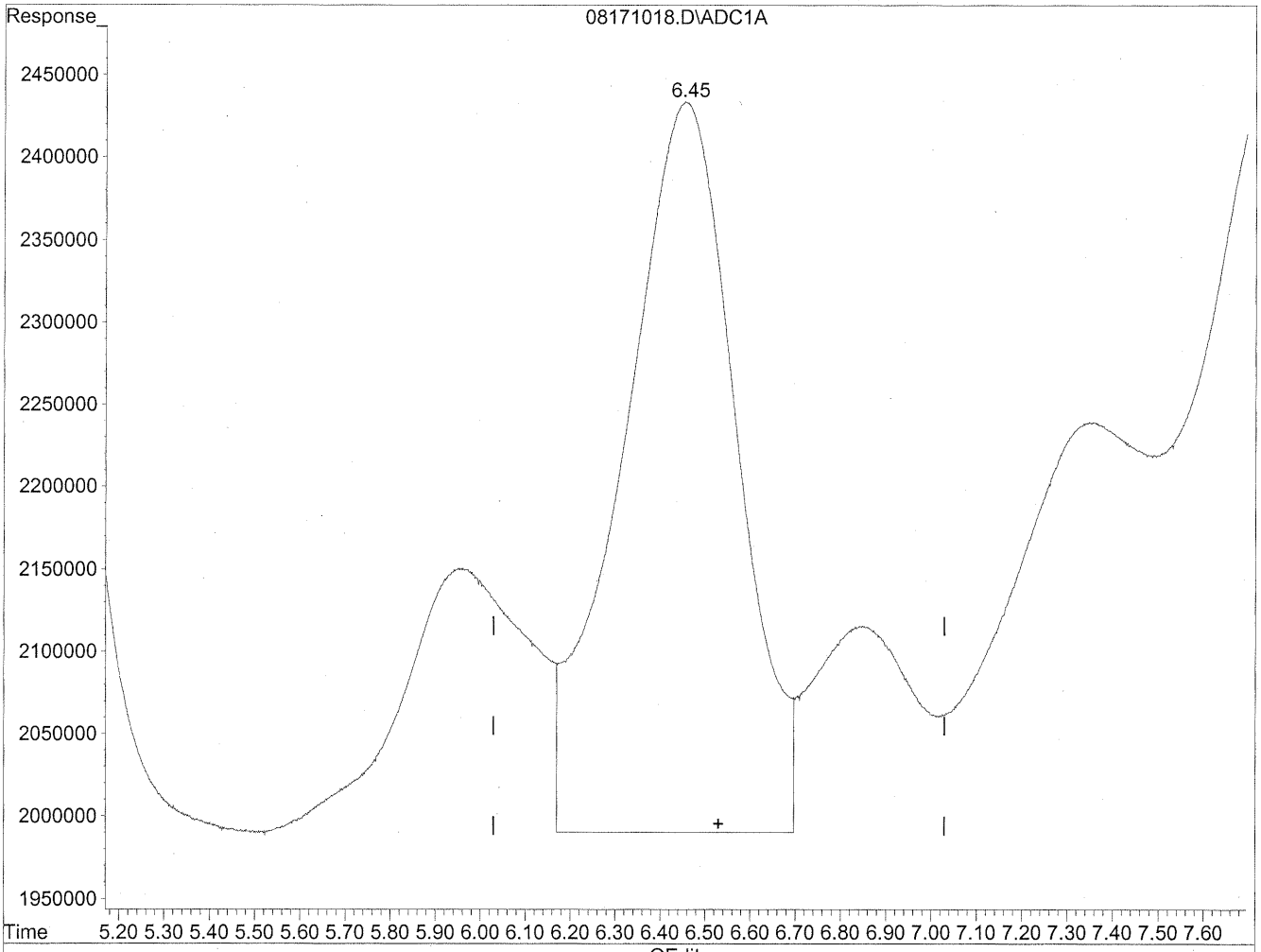
200

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171018.D Vial: 31
Acq On : 18 Aug 2009 8:09 pm Operator: HC
Sample : P0902772-008 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.46min 1182.769ng/ml

response 77908191

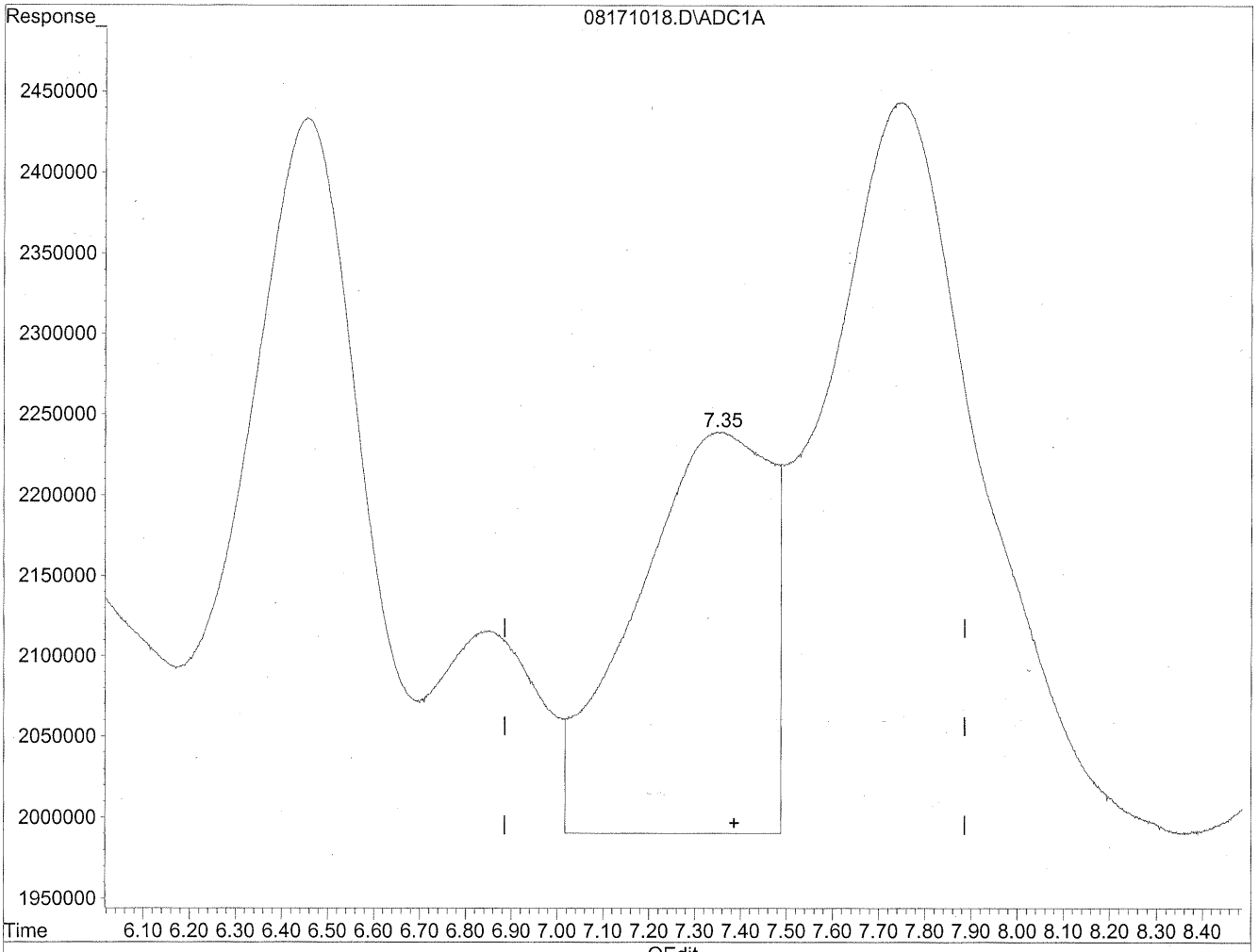
(+) = Expected Retention Time

08171018.D TO110709.M Sat Aug 22 14:34:27 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171018.D Vial: 31
Acq On : 18 Aug 2009 8:09 pm Operator: HC
Sample : P0902772-008 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.35min 650.713ng/ml

response 50918939

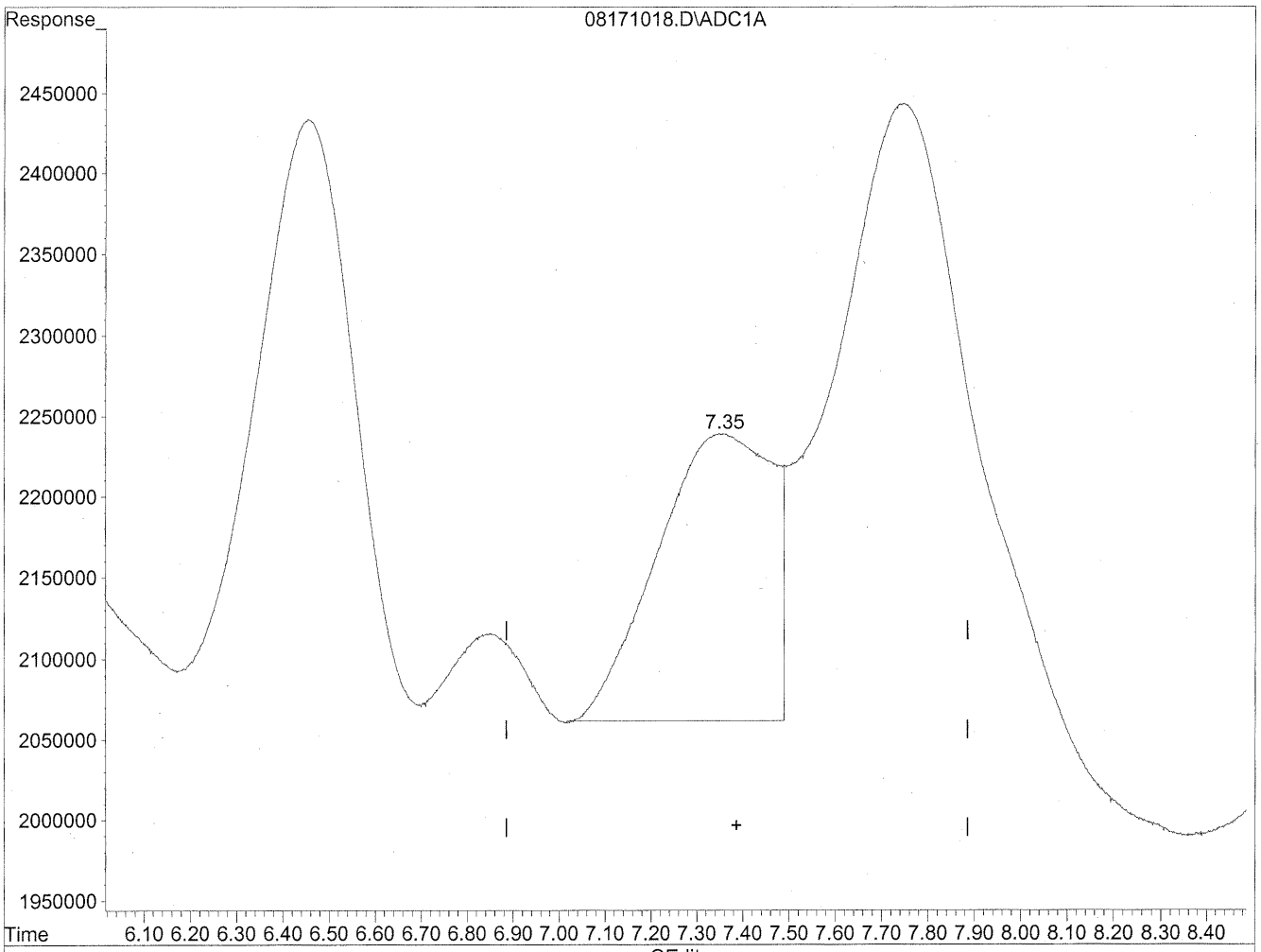
202

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171018.D Vial: 31
Acq On : 18 Aug 2009 8:09 pm Operator: HC
Sample : P0902772-008 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.35min 393.376ng/ml m

response 30782051

HC 8/22/09
RC

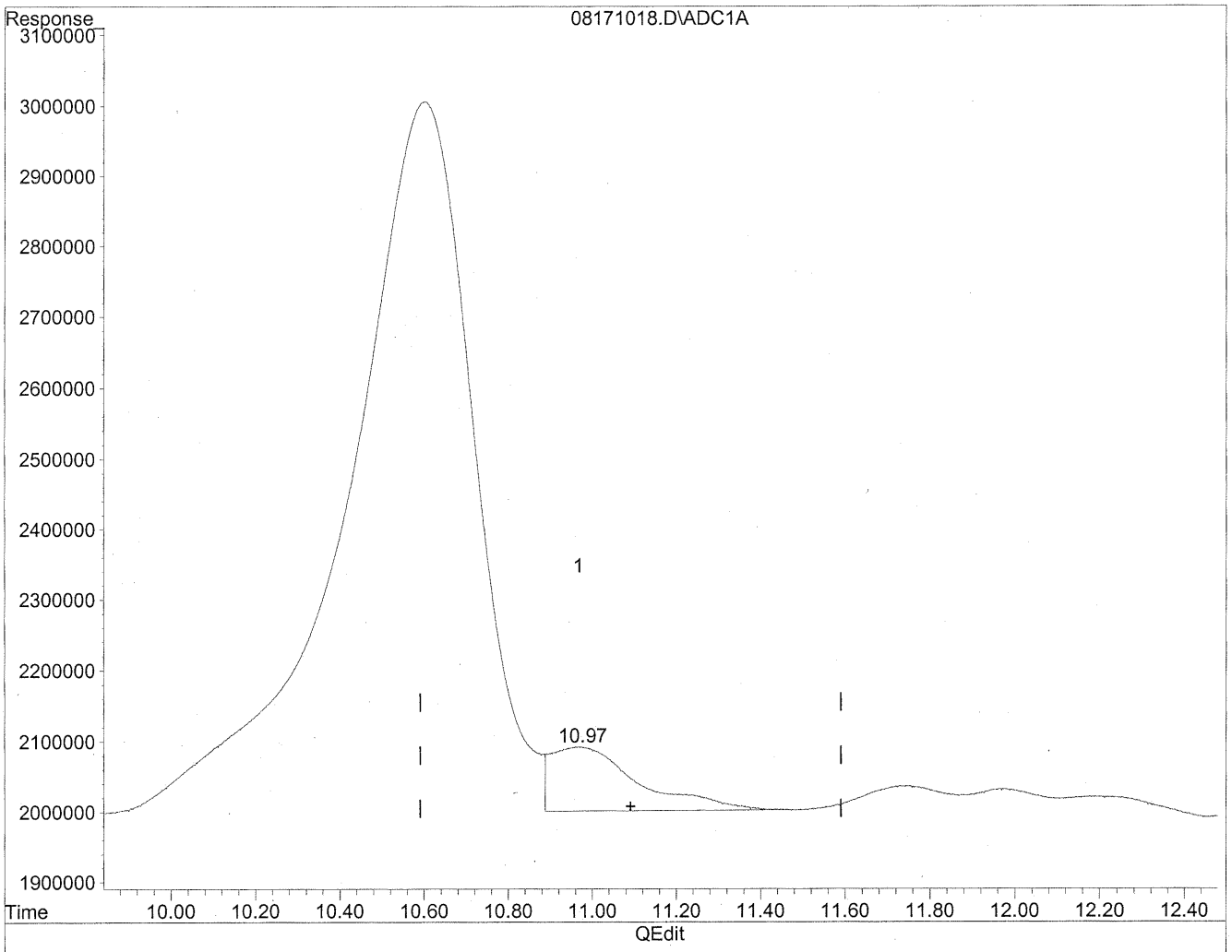
KP 8/23/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171018.D Vial: 31
Acq On : 18 Aug 2009 8:09 pm Operator: HC
Sample : P0902772-008 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.97min 267.850ng/ml

response 13128216

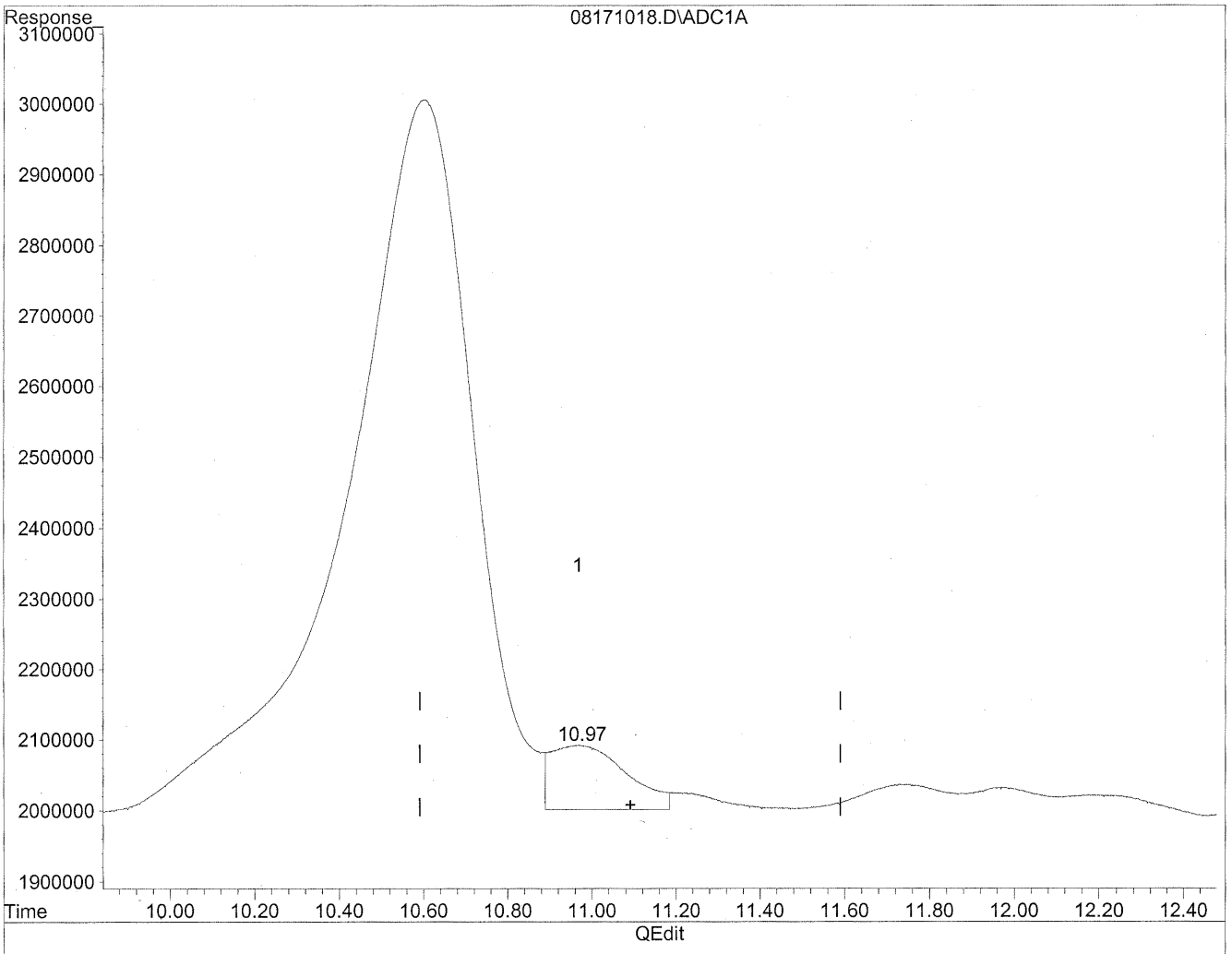
(+) = Expected Retention Time

08171018.D TO110709.M Sat Aug 22 14:35:10 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171018.D Vial: 31
Acq On : 18 Aug 2009 8:09 pm Operator: HC
Sample : P0902772-008 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.97min 230.567ng/ml m

response 11300857

HC 8/22/09
BC
KK 8/23/09

205

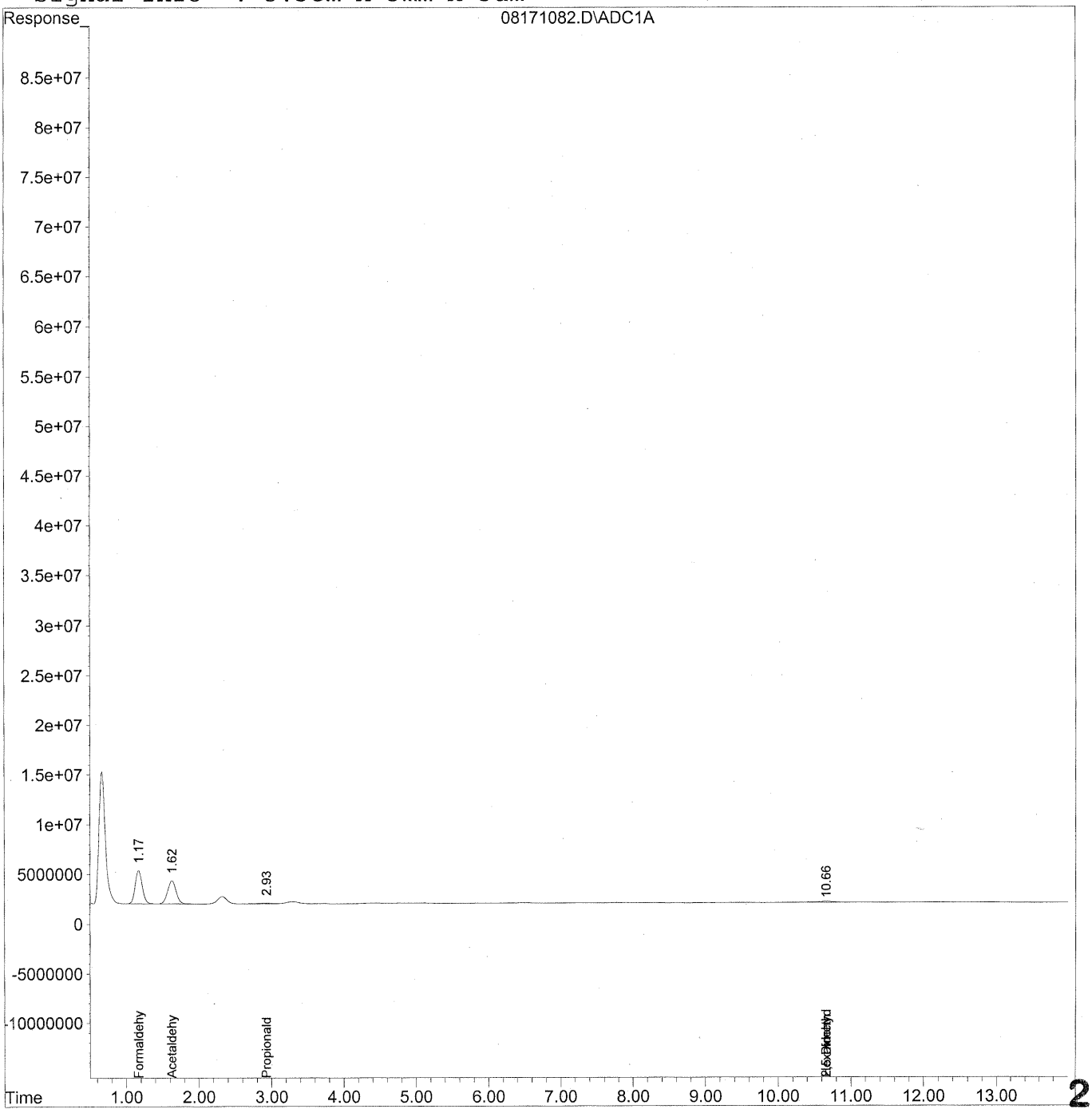
(+) = Expected Retention Time
08171018.D TO110709.M Sat Aug 22 14:35:20 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171082.D Vial: 7
Acq On : 19 Aug 2009 12:12 pm Operator: HC
Sample : P0902772-008 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:22 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



206

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171082.D Vial: 7
Acq On : 19 Aug 2009 12:12 pm Operator: HC
Sample : P0902772-008 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:22 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

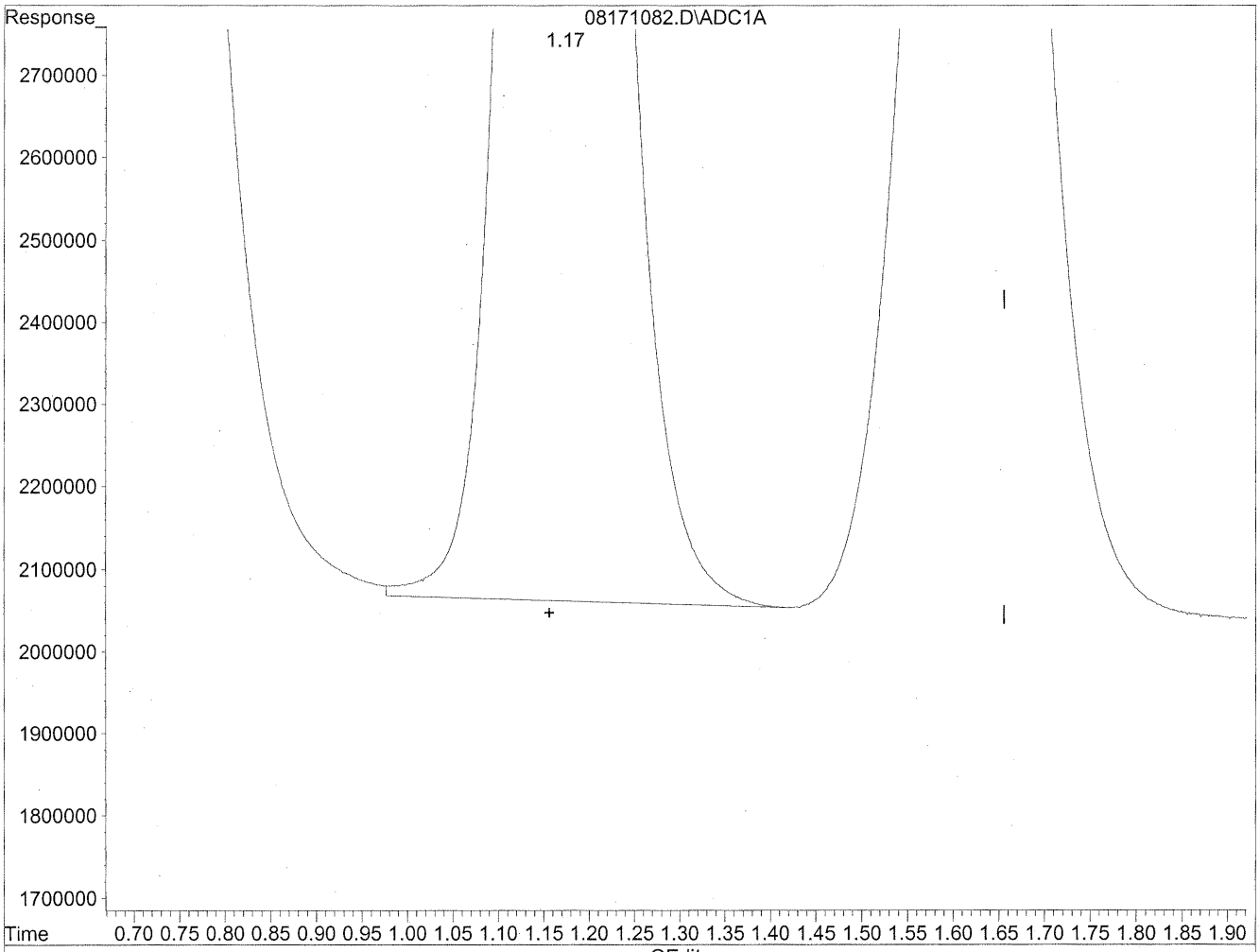
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	222269684	1210.742	ng/mlm
2)	Acetaldehyde	1.62	190051829	1355.349	ng/ml
3)	Propionaldehyde	2.93	10032483	94.029	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	10.66	18716806	277.929	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.66	18716806	381.871	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171082.D Vial: 7
Acq On : 19 Aug 2009 12:12 pm Operator: HC
Sample : P0902772-008 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 1218.655ng/ml

response 223722366

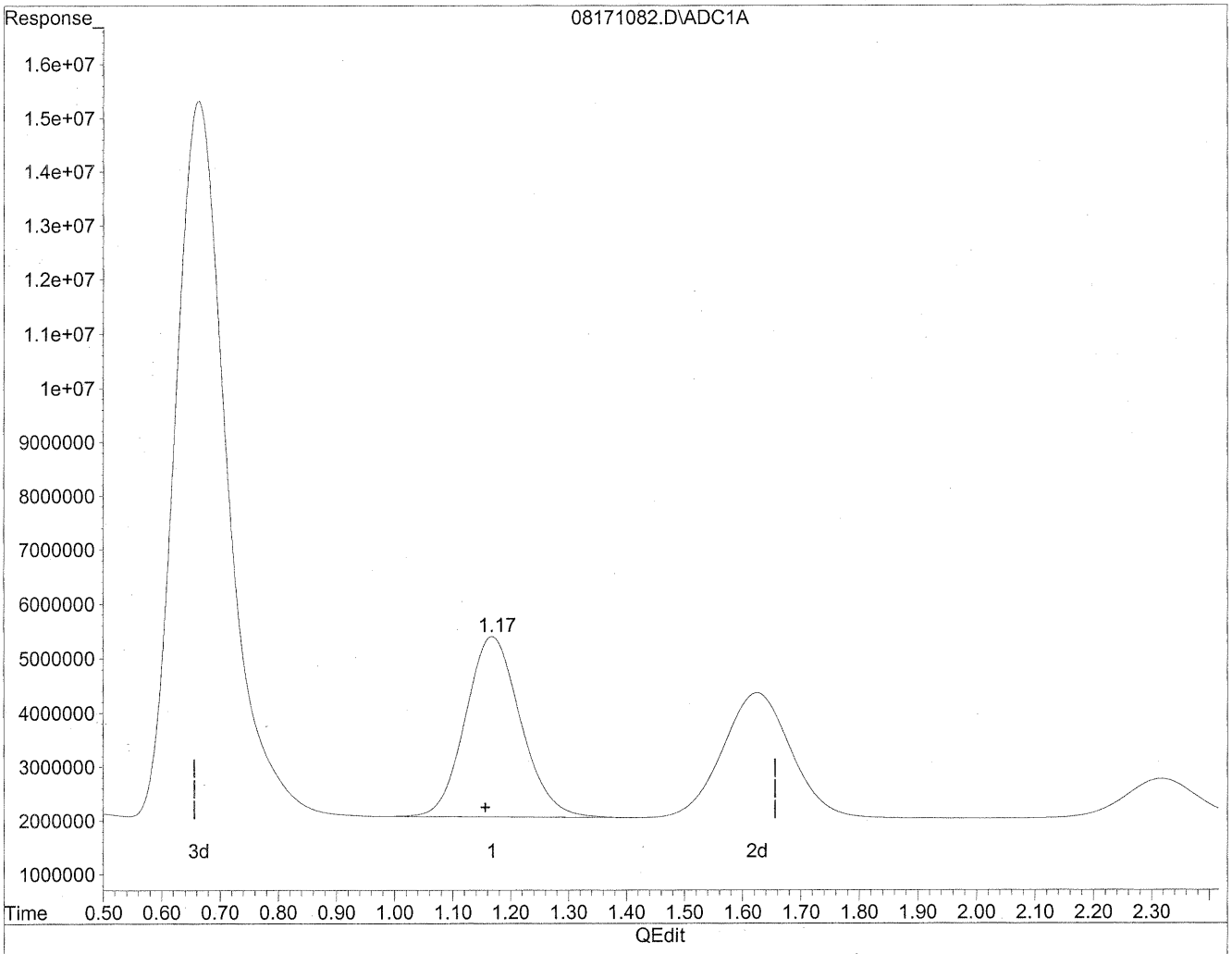
208

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171082.D Vial: 7
Acq On : 19 Aug 2009 12:12 pm Operator: HC
Sample : P0902772-008 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 1210.742ng/ml m

response 222269684

HC 8/22/09
LC
KP 8/23/09

209

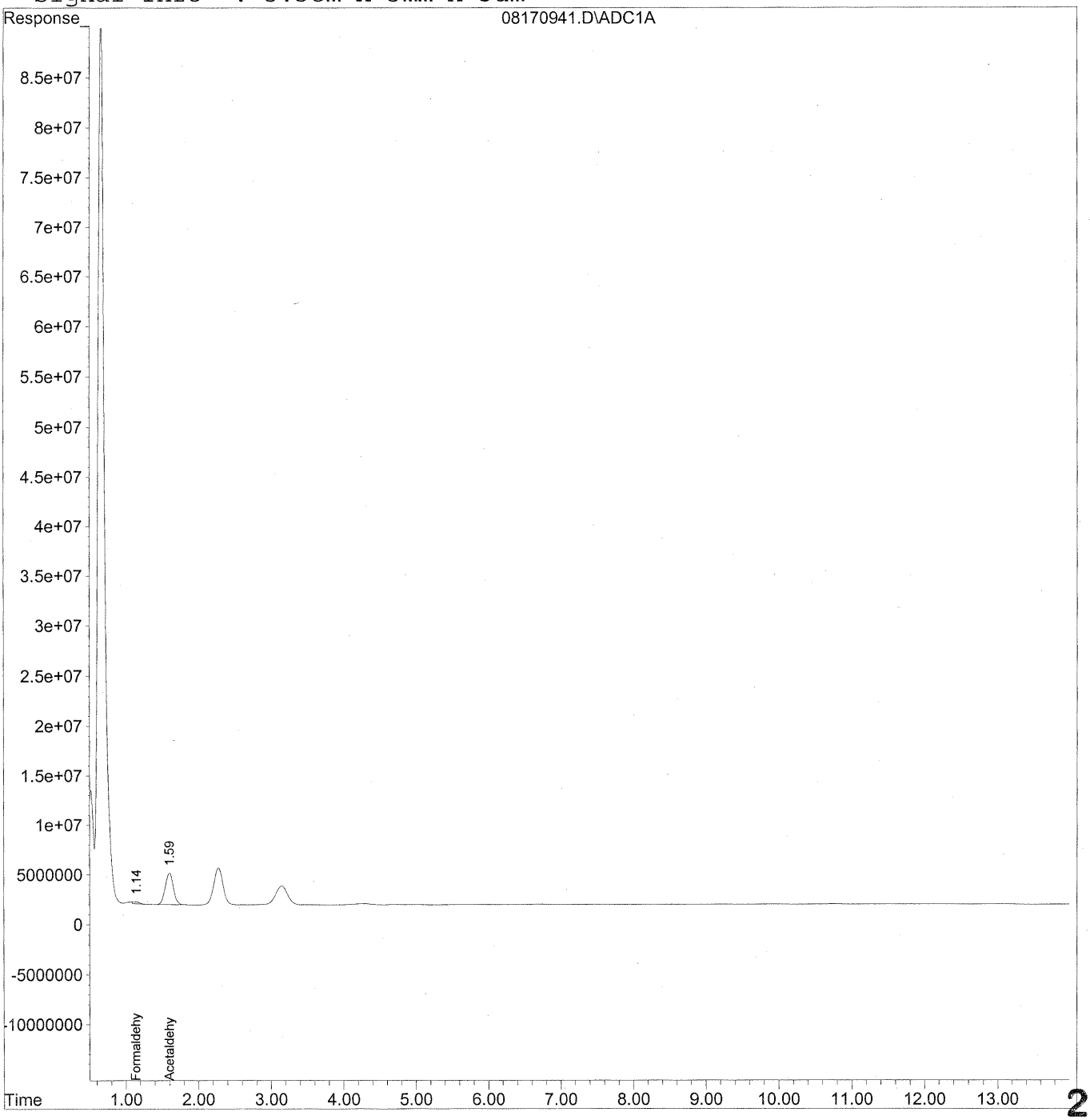
(+) = Expected Retention Time
08171082.D TO110709.M Sat Aug 22 17:22:48 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170941.D Vial: 40
Acq On : 18 Aug 2009 12:52 am Operator: HC
Sample : P0902772-008 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:20 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



210

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170941.D Vial: 40
Acq On : 18 Aug 2009 12:52 am Operator: HC
Sample : P0902772-008 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:20 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

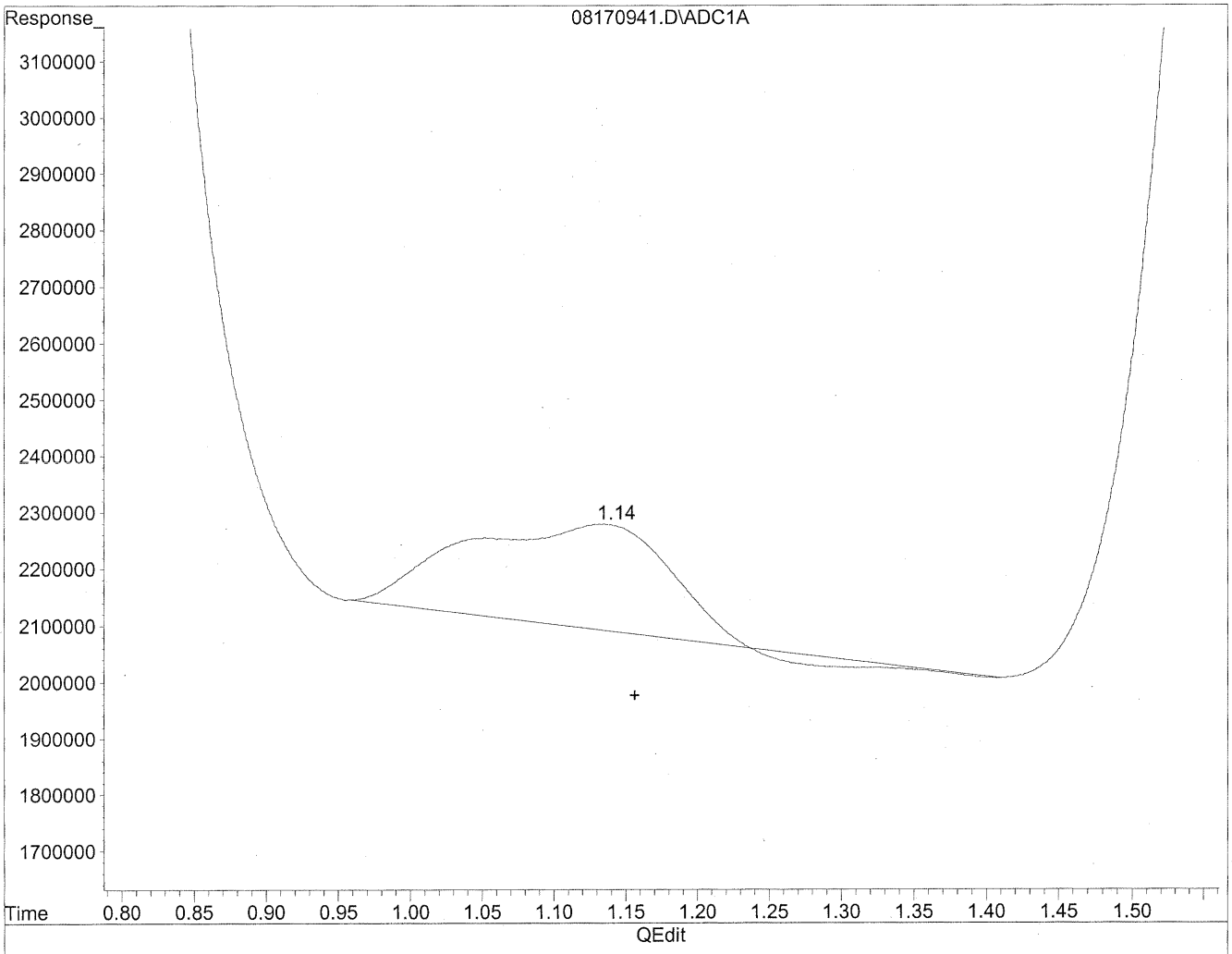
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.14	11998752	65.359	ng/mlm
2)	Acetaldehyde	1.59	248250108	1770.389	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170941.D Vial: 40
Acq On : 18 Aug 2009 12:52 am Operator: HC
Sample : P0902772-008 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.13min 94.452ng/ml

response 17339701

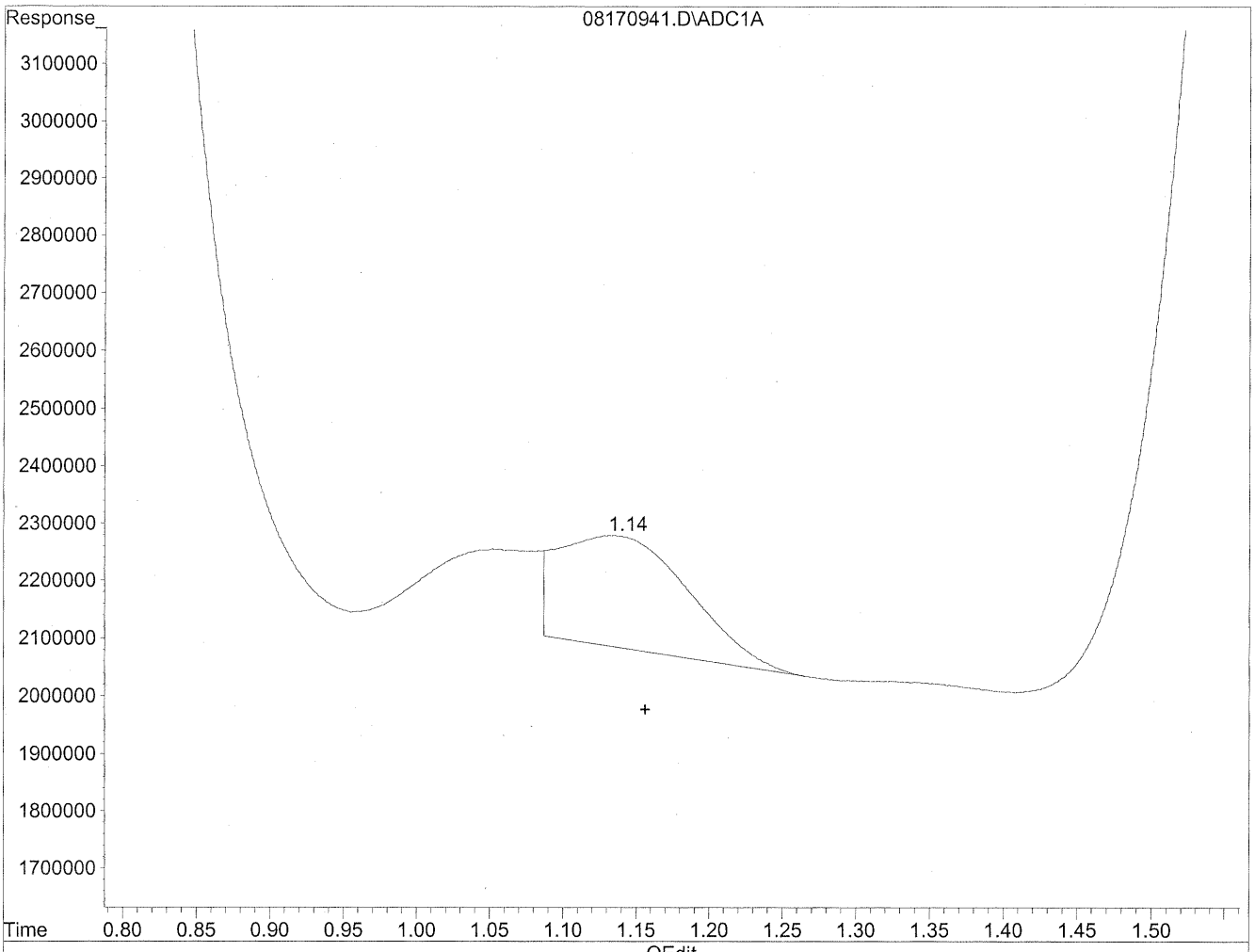
212

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170941.D Vial: 40
Acq On : 18 Aug 2009 12:52 am Operator: HC
Sample : P0902772-008 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.14min 65.359ng/ml m

response 11998752

HC 8/21/09 SP

KR 8/23/09

213

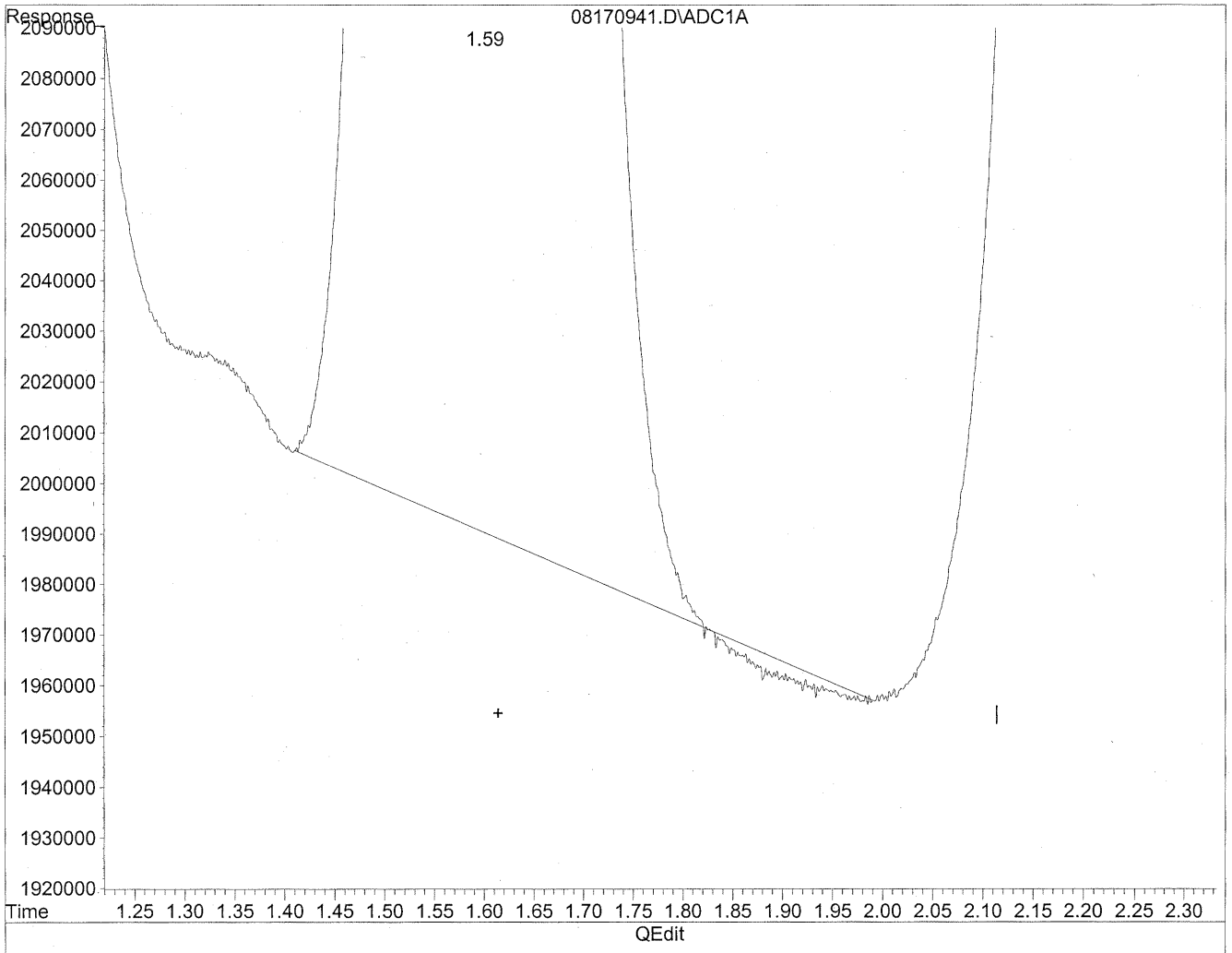
(+) = Expected Retention Time

08170941.D TO110709.M Fri Aug 21 18:19:48 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170941.D Vial: 40
Acq On : 18 Aug 2009 12:52 am Operator: HC
Sample : P0902772-008 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.59min 1768.259ng/ml

response 247951494

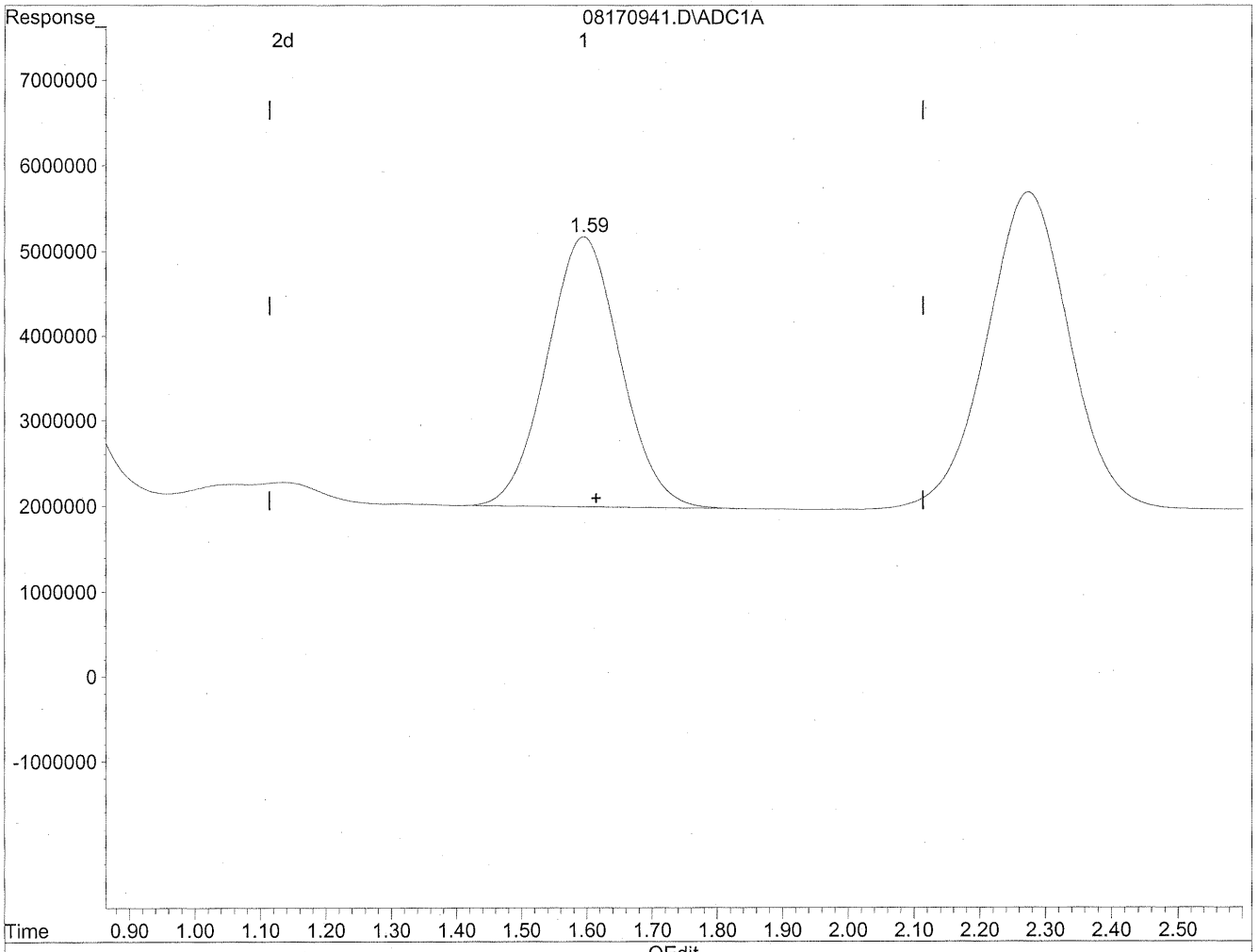
214

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170941.D Vial: 40
Acq On : 18 Aug 2009 12:52 am Operator: HC
Sample : P0902772-008 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.59min 1770.389ng/ml m

response 248250108

HC 8/21/09

KP 8/23/09

215

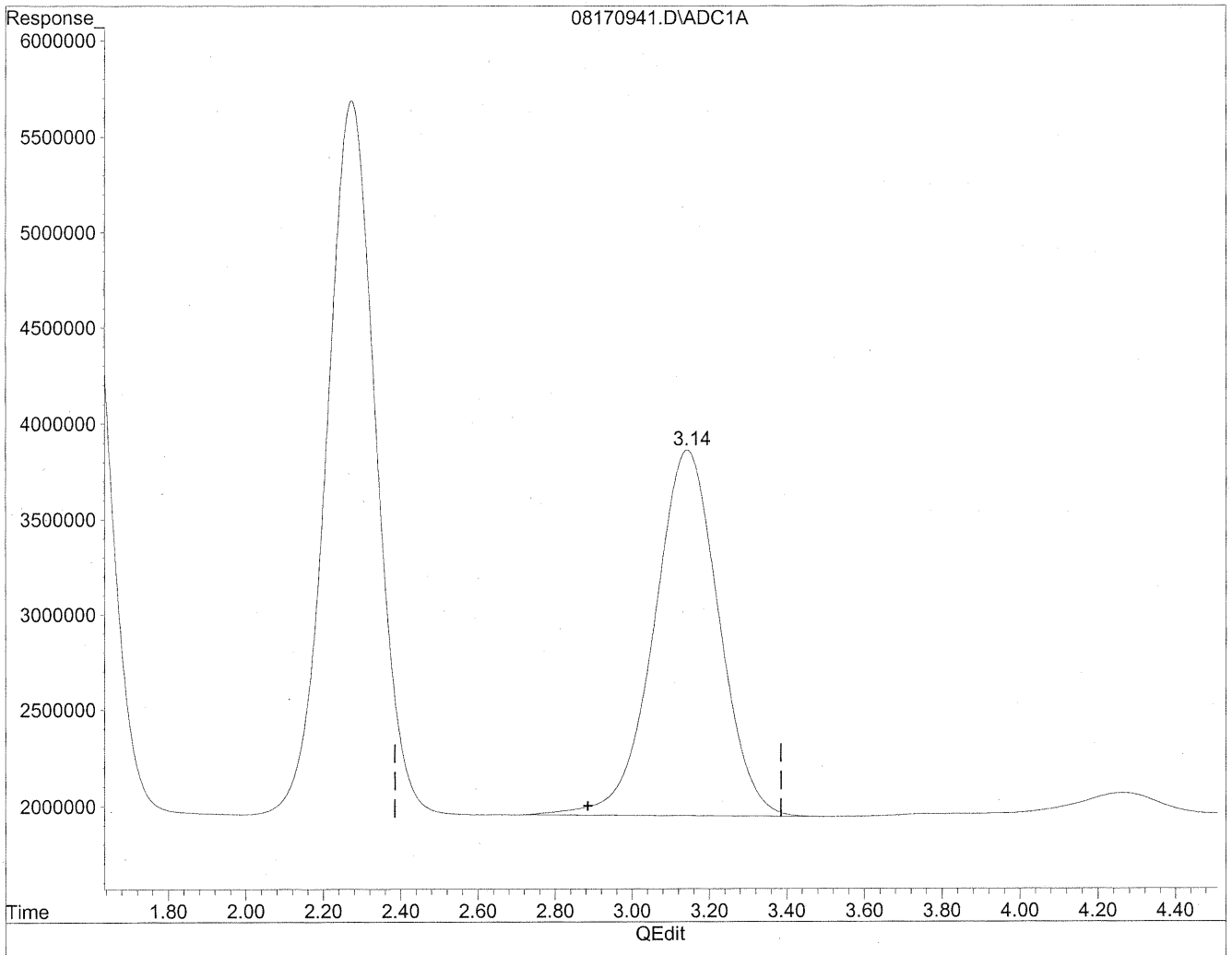
(+) = Expected Retention Time

08170941.D TO110709.M Fri Aug 21 18:20:04 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170941.D Vial: 40
Acq On : 18 Aug 2009 12:52 am Operator: HC
Sample : P0902772-008 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.14min 2083.726ng/ml

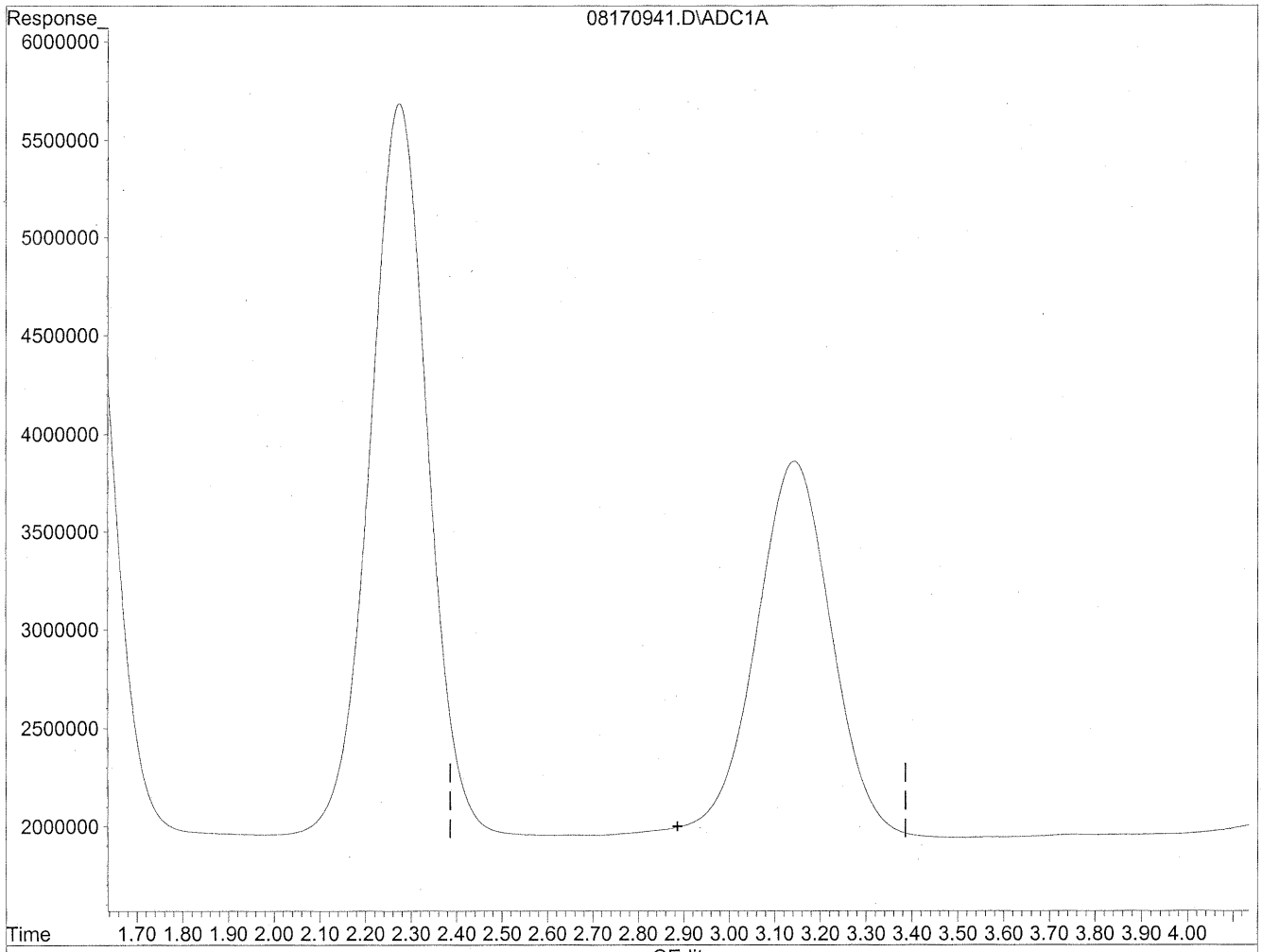
response 222323621

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170941.D Vial: 40
Acq On : 18 Aug 2009 12:52 am Operator: HC
Sample : P0902772-008 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/21/09 WP

WP 8/23/09

217

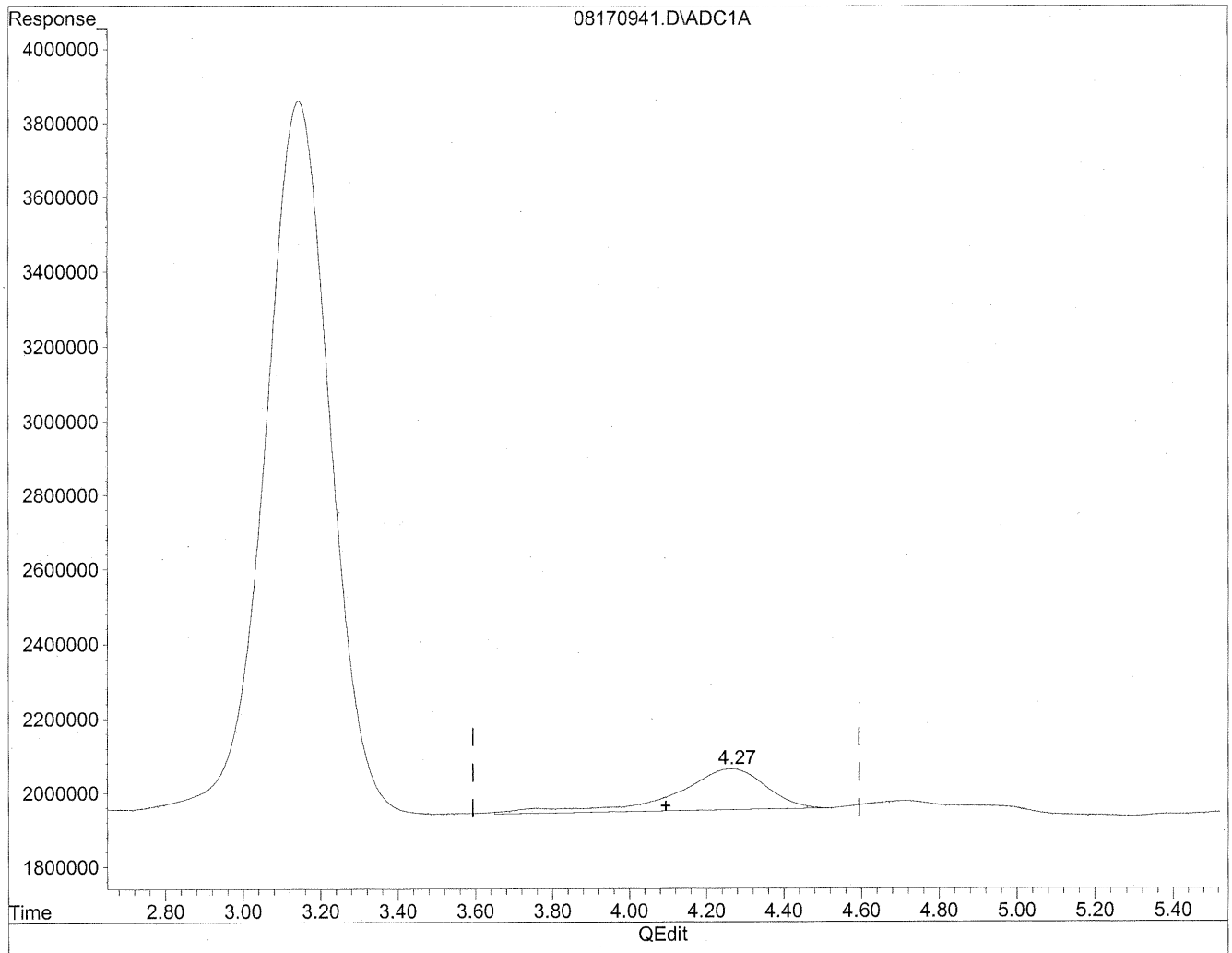
(+) = Expected Retention Time

08170941.D TO110709.M Fri Aug 21 18:20:11 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170941.D Vial: 40
Acq On : 18 Aug 2009 12:52 am Operator: HC
Sample : P0902772-008 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.26min 189.189ng/ml

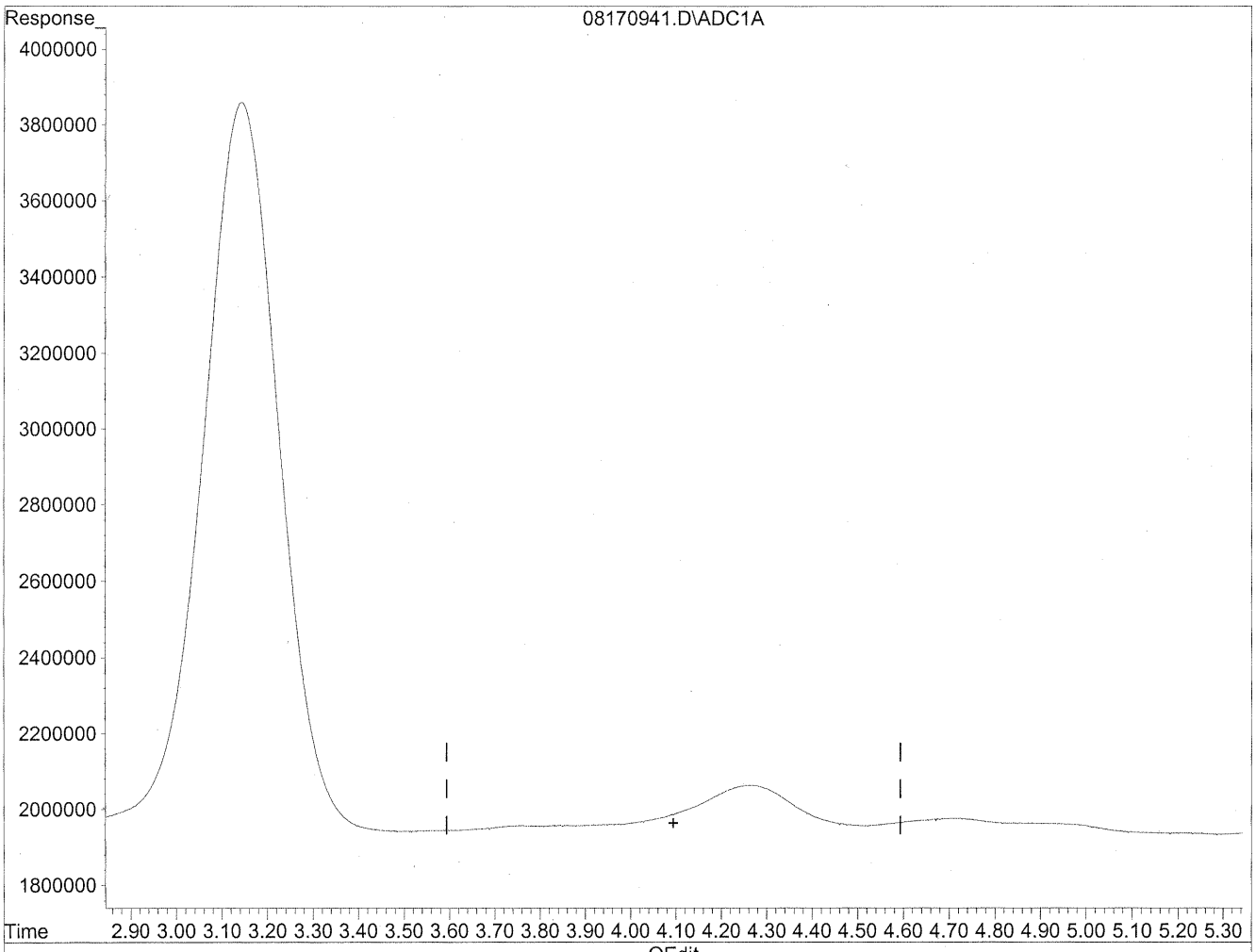
response 18429921

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170941.D Vial: 40
Acq On : 18 Aug 2009 12:52 am Operator: HC
Sample : P0902772-008 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:19 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/21/09 WP

KP 8/23/09

(+) = Expected Retention Time

08170941.D TO110709.M Fri Aug 21 18:20:17 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **100454**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P0902772-009

Test Code: EPA Method TO-11A
Instrument ID: Waters LC Module I Plus/UV_Vis 360/LC1
Analyst: Hani Cherazaie
Sampling Media: Silica Gel DNPH Tube
Test Notes: **BC**

Date Collected: 8/10/09
Date Received: 8/12/09
Date Analyzed: 8/18/09
Desorption Volume: 1.0 ml
Volume Sampled: NA Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	Formaldehyde	< 100	NA	NA	NA	NA	
75-07-0	Acetaldehyde	< 100	NA	NA	NA	NA	
123-38-6	Propionaldehyde	< 100	NA	NA	NA	NA	
4170-30-3	Crotonaldehyde, Total	< 100	NA	NA	NA	NA	
123-72-8	Butyraldehyde	< 100	NA	NA	NA	NA	
100-52-7	Benzaldehyde	< 100	NA	NA	NA	NA	
590-86-3	Isovaleraldehyde	< 100	NA	NA	NA	NA	
110-62-3	Valeraldehyde	< 100	NA	NA	NA	NA	
529-20-4	o-Tolualdehyde	< 100	NA	NA	NA	NA	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	NA	NA	NA	NA	
66-25-1	n-Hexaldehyde	< 100	NA	NA	NA	NA	
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	NA	NA	NA	NA	

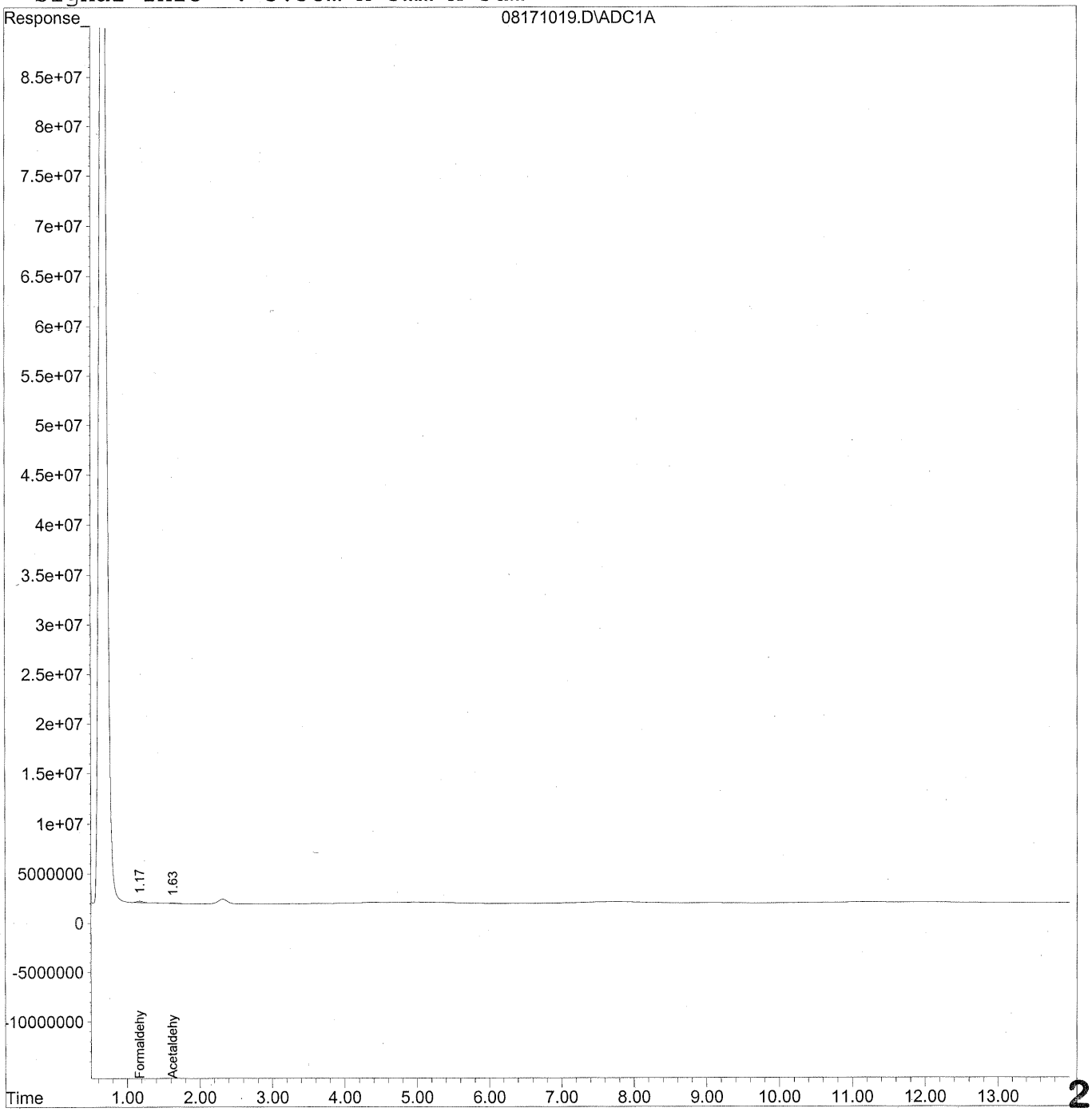
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.

Verified By: _____ Date: 8/25/09

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:37 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



221

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:37 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

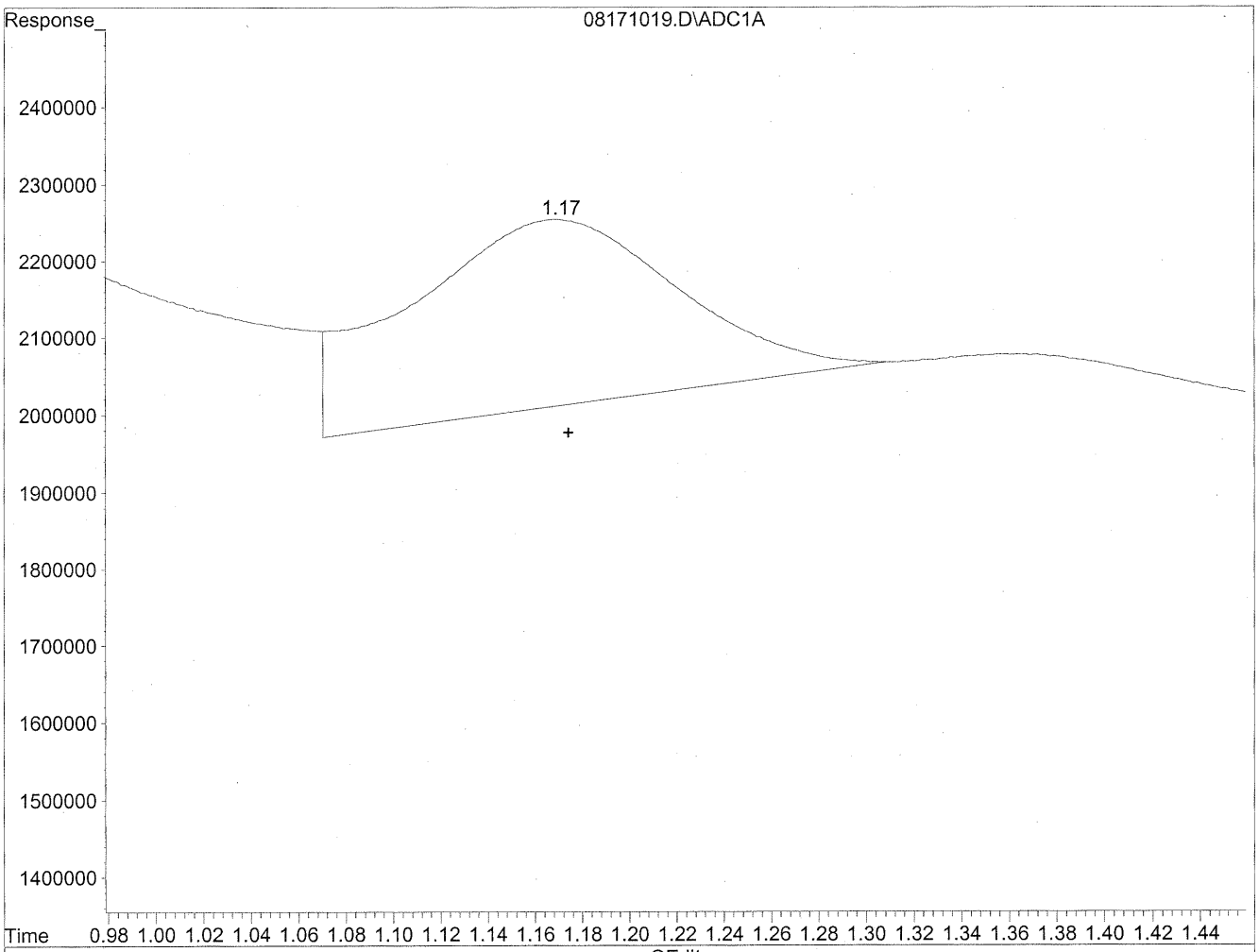
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	9645408	52.540	ng/mlm
2)	Acetaldehyde	1.63	3068771	21.885	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	0.00	0	N.D.	ng/mld
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/mld
8)	Valeraldehyde	0.00	0	N.D.	ng/mld
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/mld

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.17min 106.394ng/ml

response 19532005

223

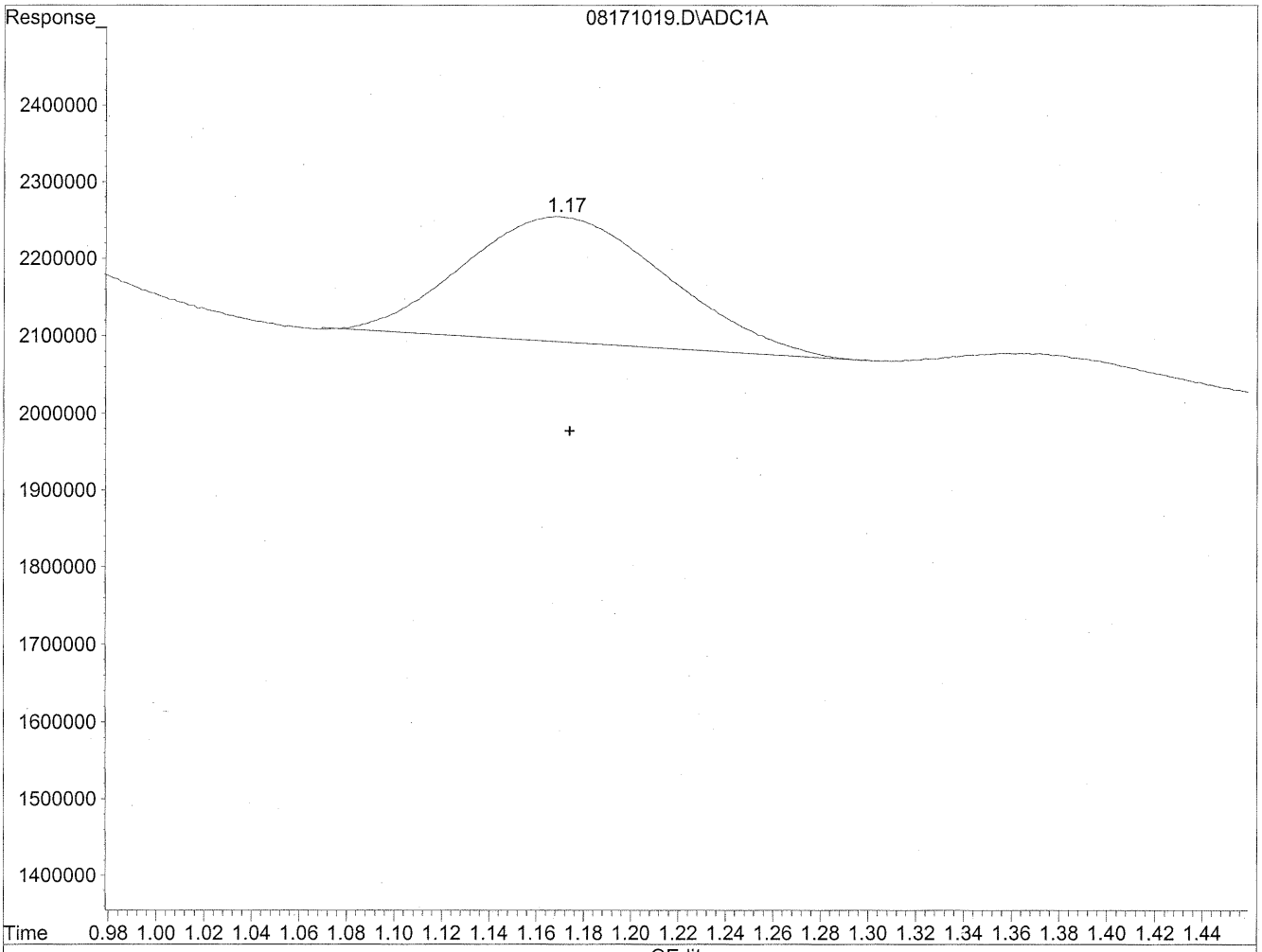
(+) = Expected Retention Time

08171019.D TO110709.M Sat Aug 22 14:37:05 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.17min 52.540ng/ml m

response 9645408

tc 8/22/09

1c

KP 8/23/09

224

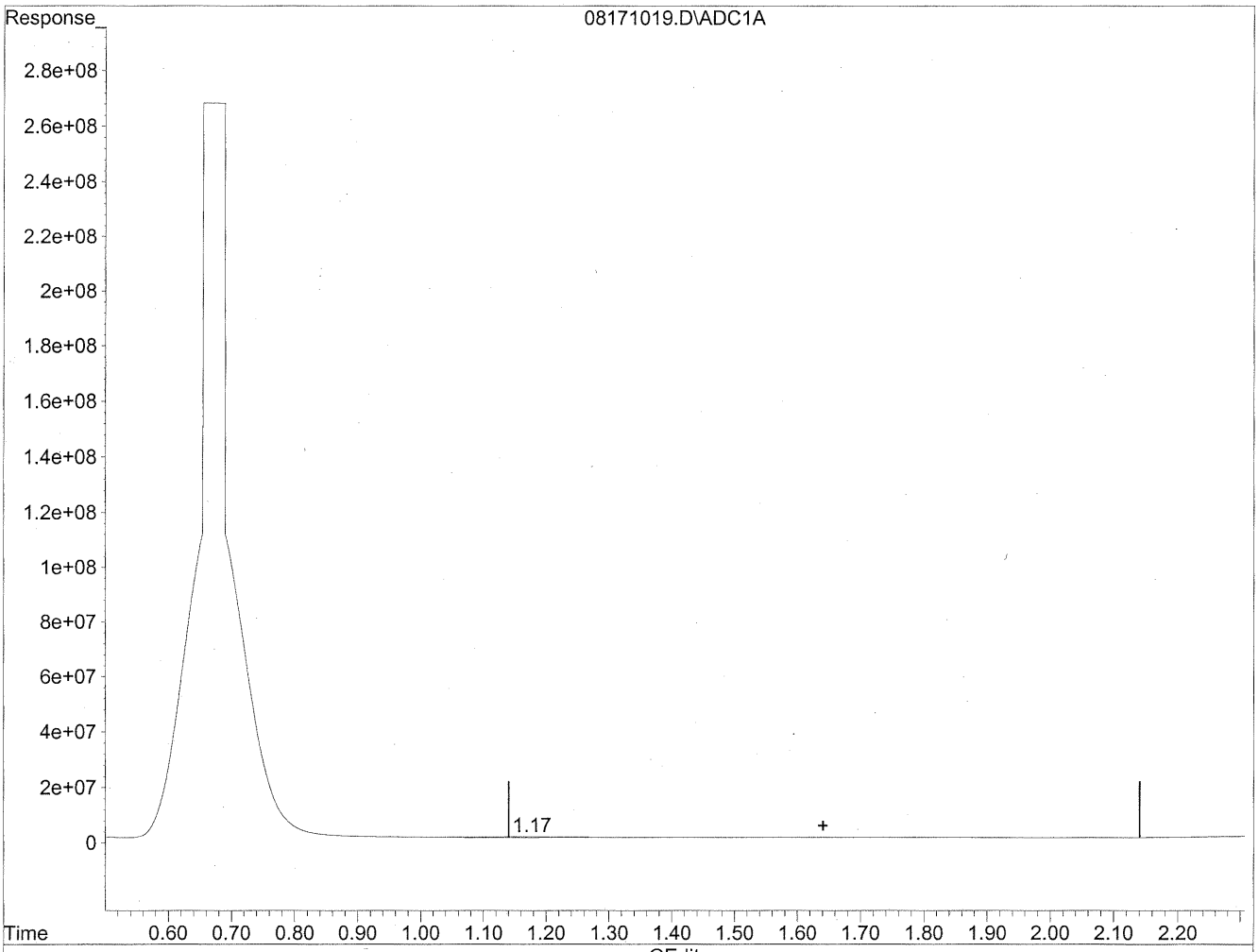
(+) = Expected Retention Time

08171019.D TO110709.M Sat Aug 22 14:37:11 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.17min 139.292ng/ml

response 19532005

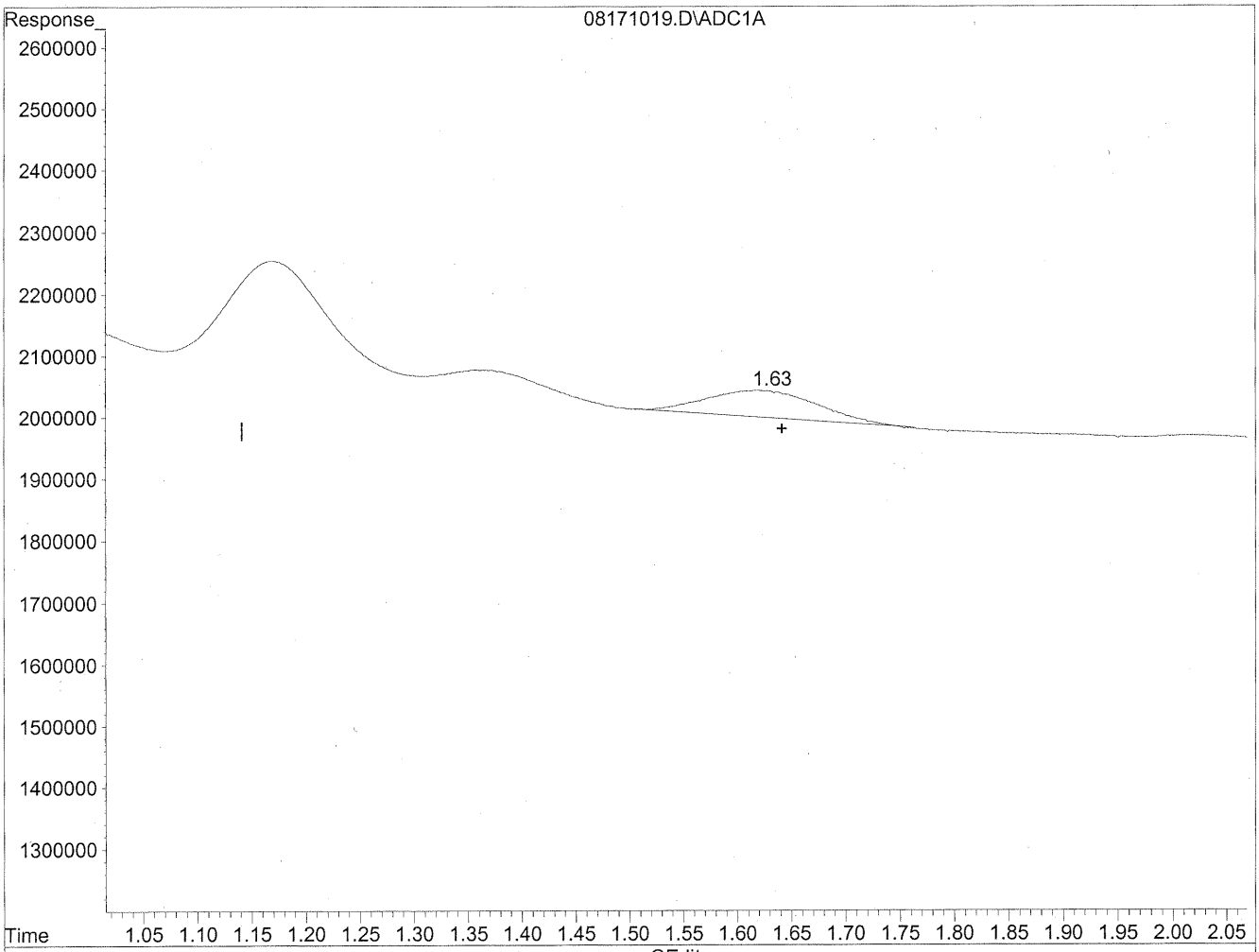
225

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.63min 21.885ng/ml m

response 3068771

HL 8/22/09 IC

KB 8/23/09

226

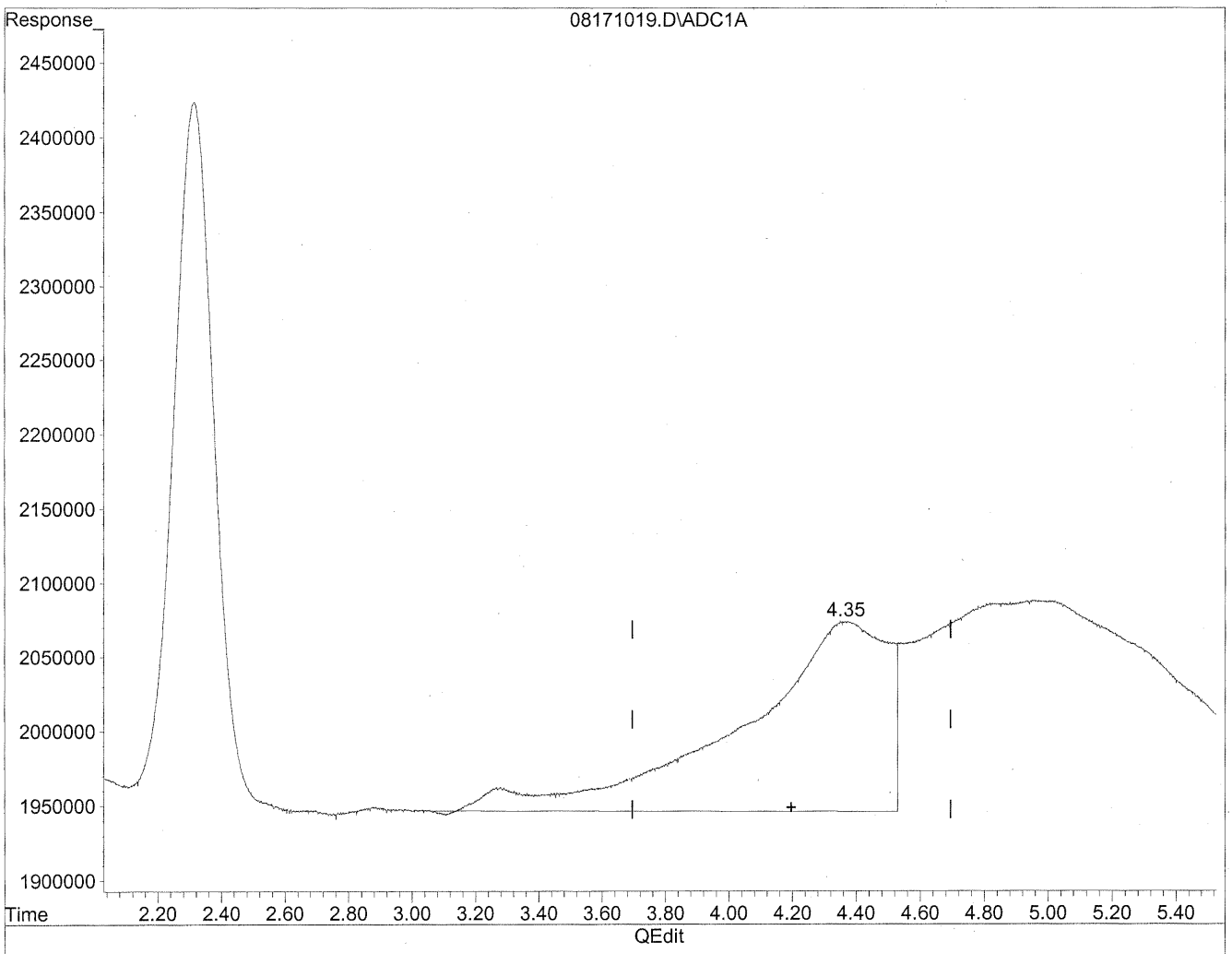
(+) = Expected Retention Time

08171019.D TO110709.M Sat Aug 22 14:37:24 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.37min 419.413ng/ml

response 40857177

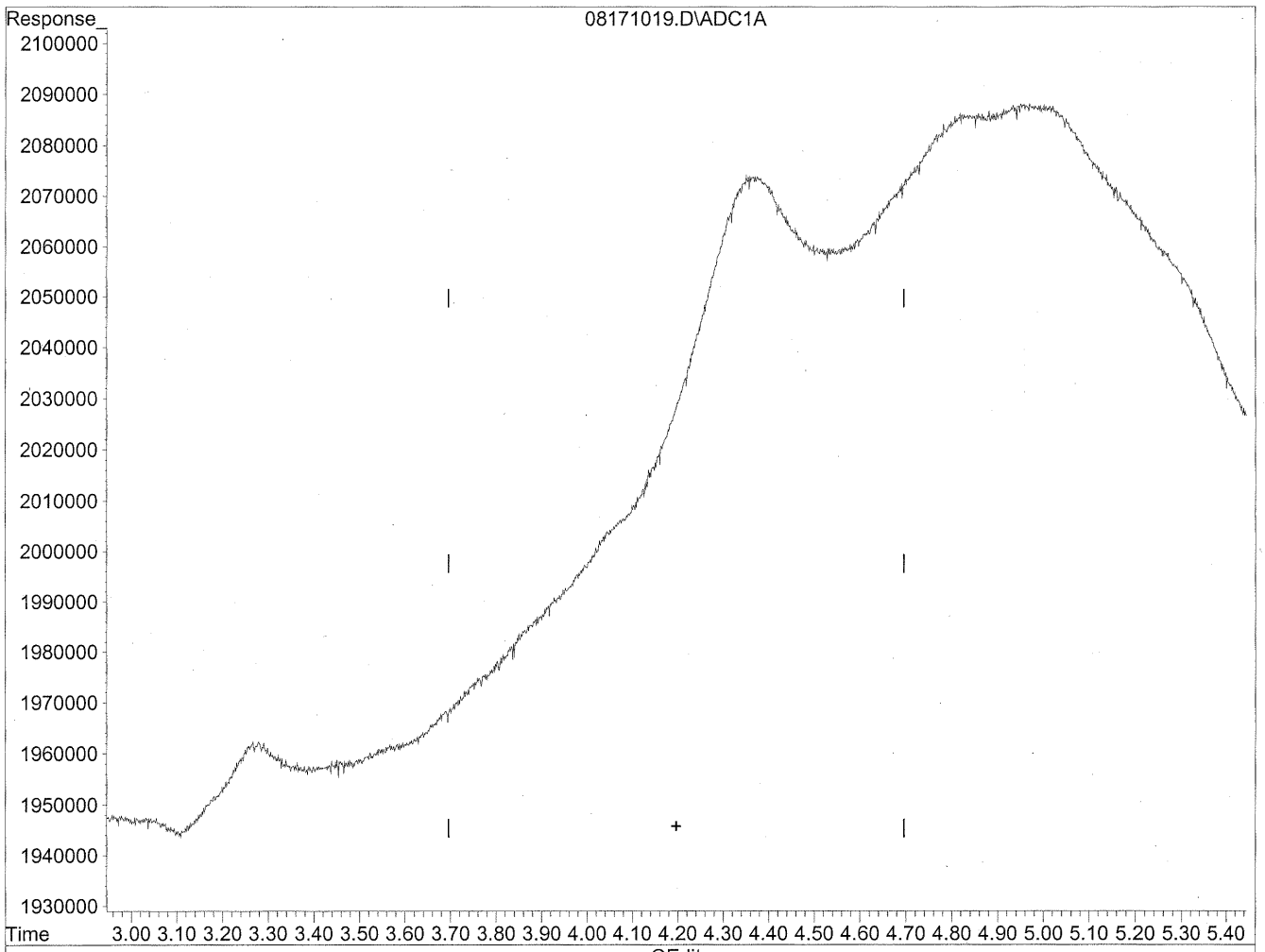
(+) = Expected Retention Time

08171019.D TO110709.M Sat Aug 22 14:37:28 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09 not reqd KR 8/23/09

228

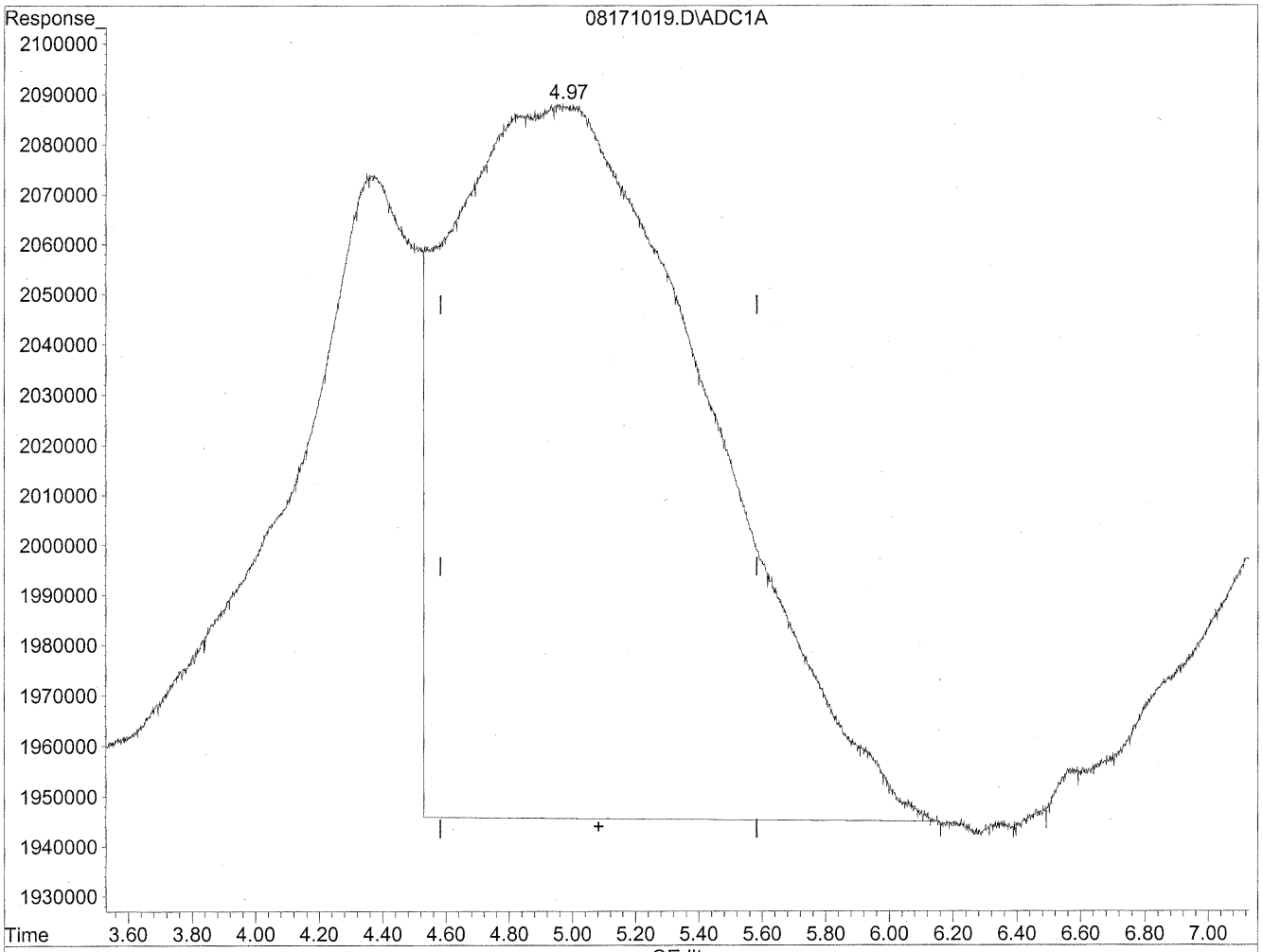
(+) = Expected Retention Time

08171019.D TO110709.M Sat Aug 22 14:37:31 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.97min 912.868ng/ml

response 80639230

229

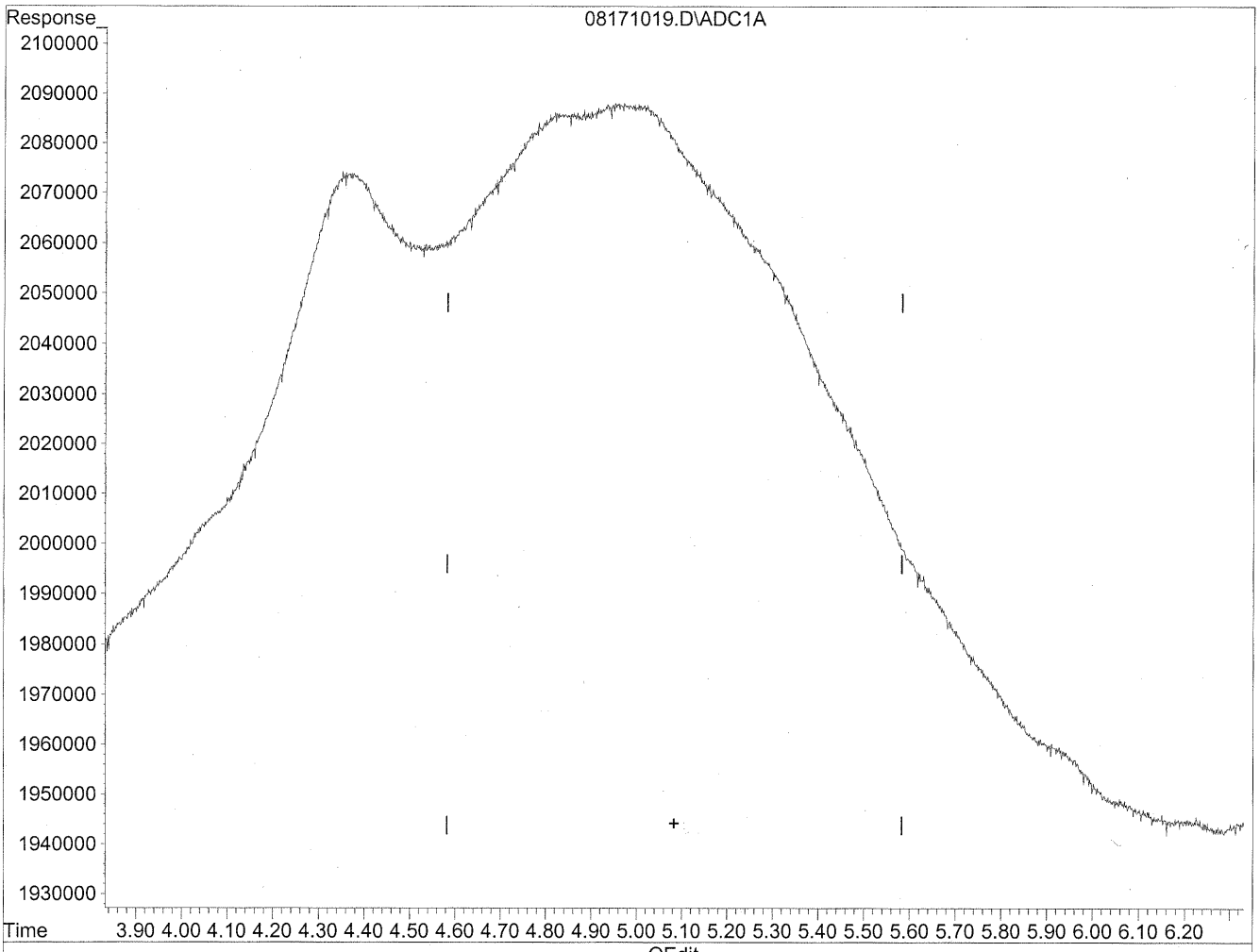
(+) = Expected Retention Time

08171019.D TO110709.M Sat Aug 22 14:37:34 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(5) Butyraldehyde

0.00min 0.000ng/ml d

response 0

tle 8/21/09 not read

KR 8/23/09

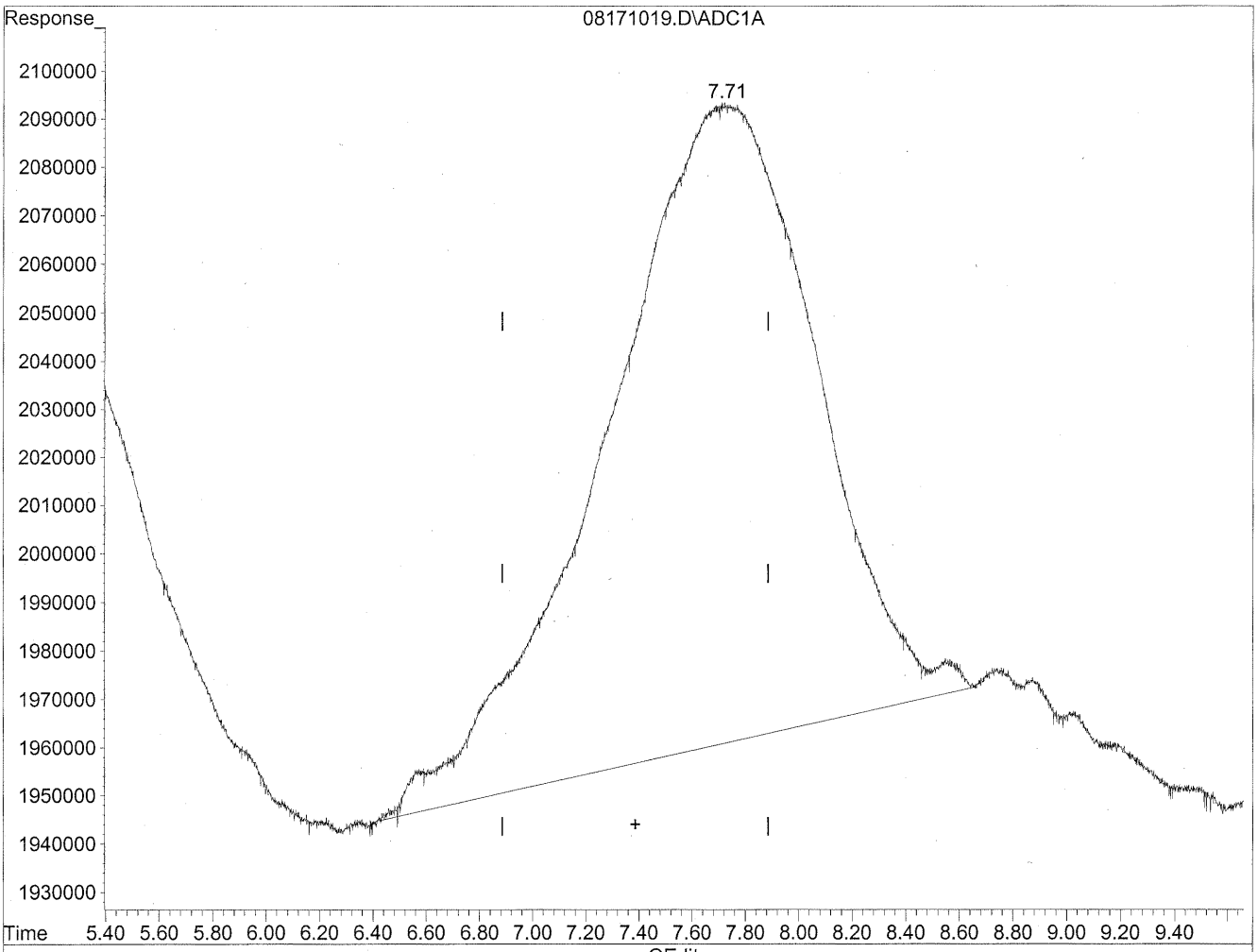
230

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.73min 926.360ng/ml

response 72488598

231

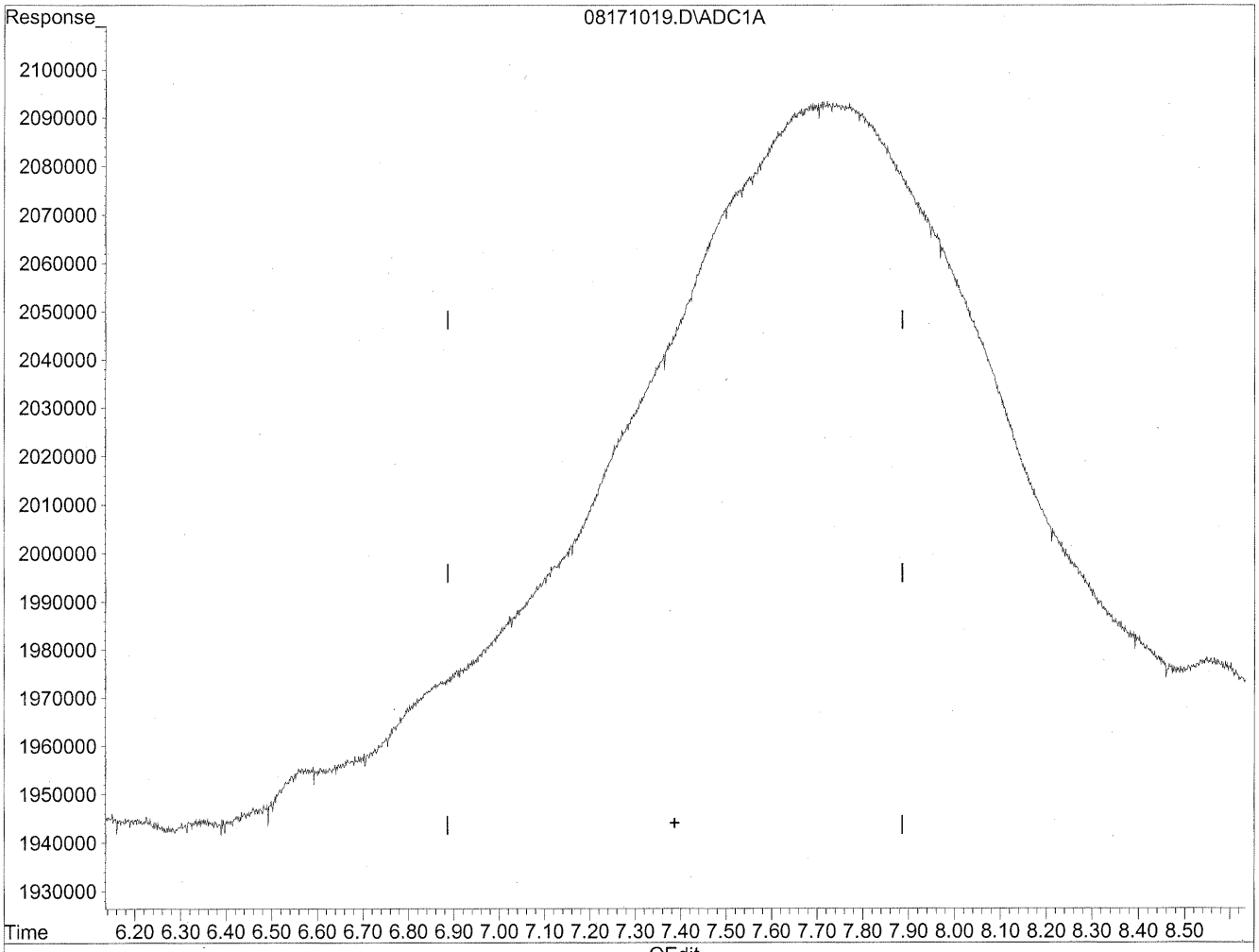
(+) = Expected Retention Time

08171019.D TO110709.M Sat Aug 22 14:37:40 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

0.00min 0.000ng/ml d

response 0

tlc 8/22/09 not read

KR 8/23/09

232

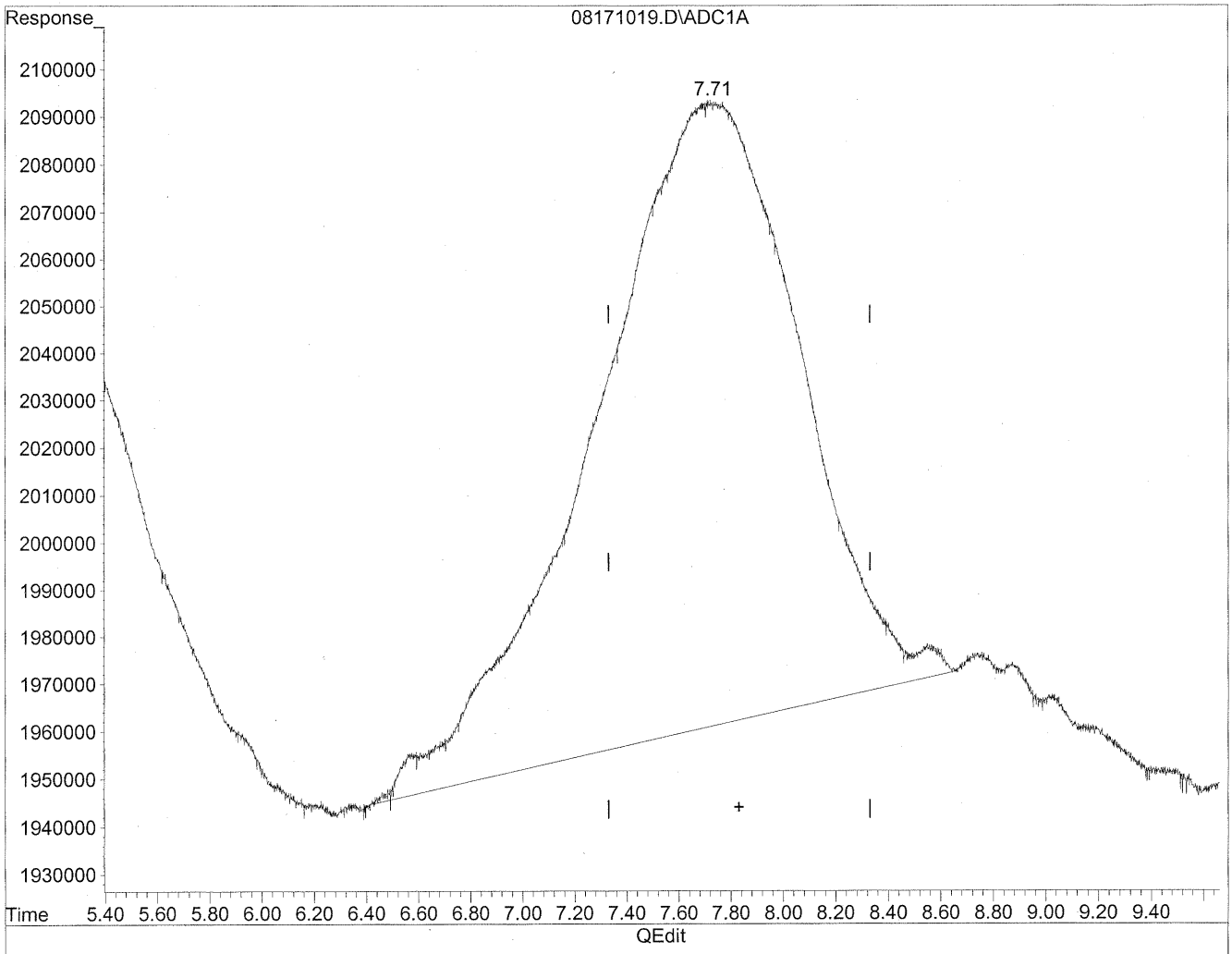
(+) = Expected Retention Time

08171019.D TO110709.M Sat Aug 22 14:37:43 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.73min 986.172ng/ml

response 72488598

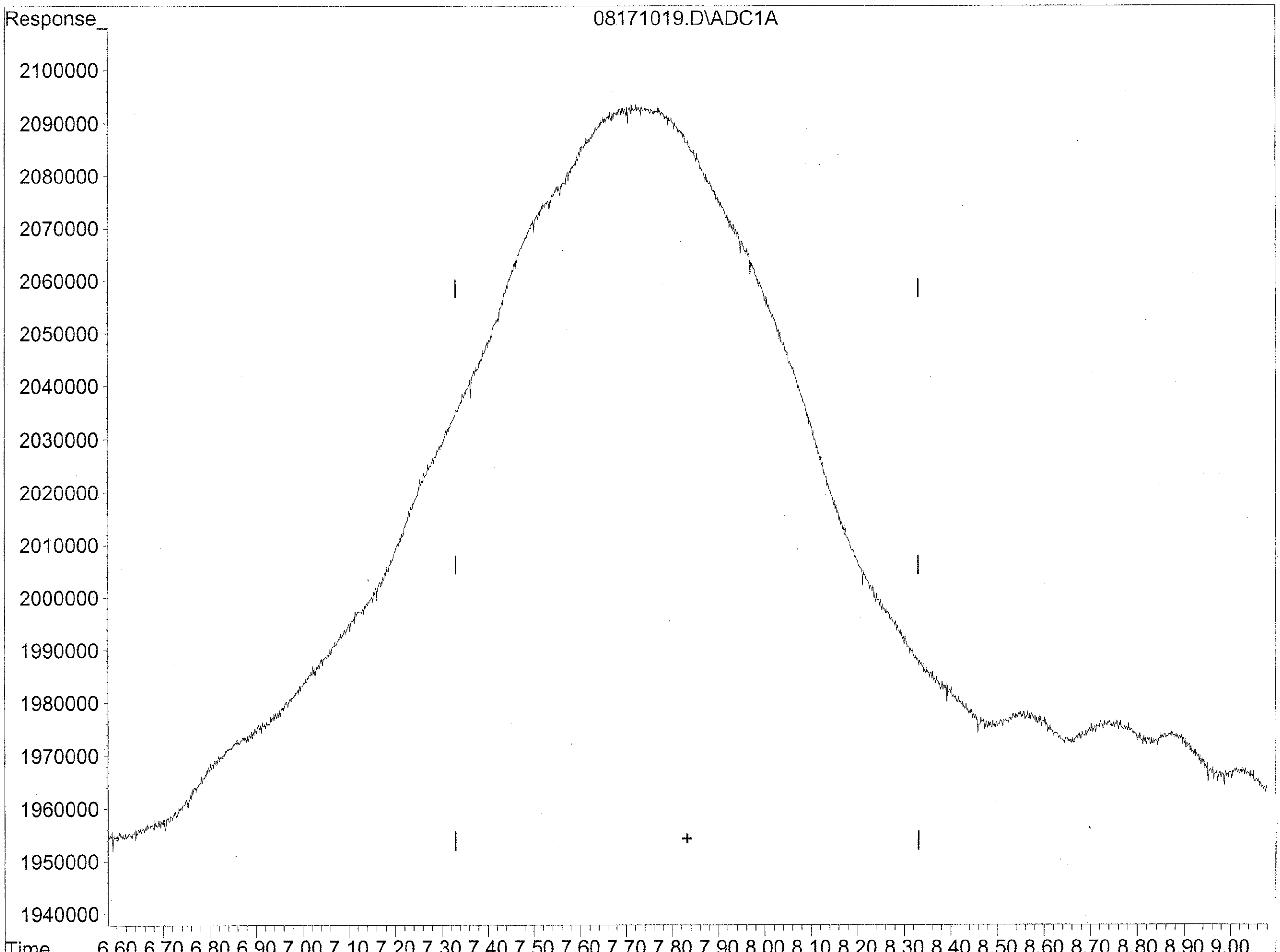
(+) = Expected Retention Time

08171019.D TO110709.M Sat Aug 22 14:37:46 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(8) Valeraldehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09 not real

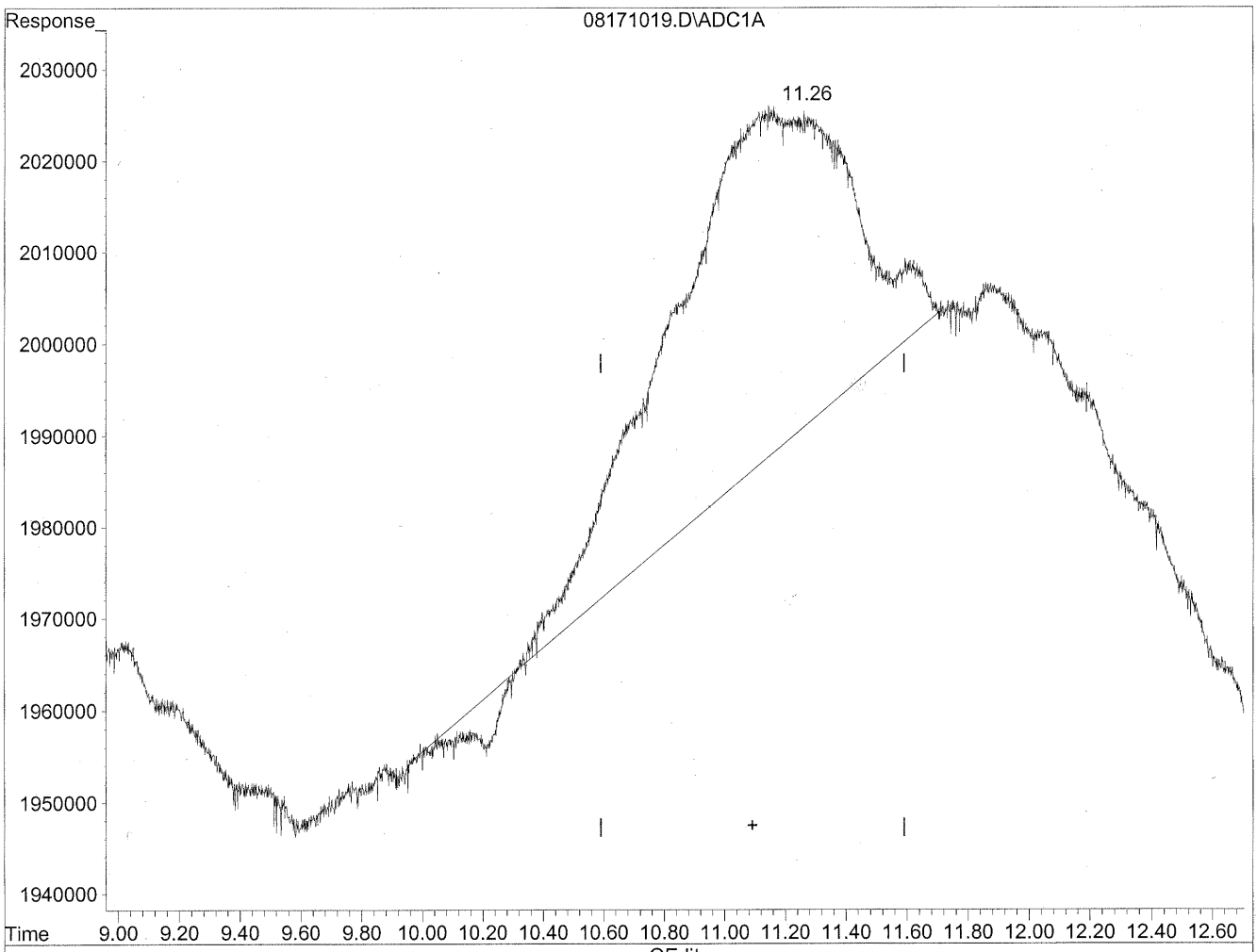
KR 8/23/09

(+) = Expected Retention Time
08171019.D TO110709.M Sat Aug 22 14:37:49 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

11.15min 316.397ng/ml

response 15507693

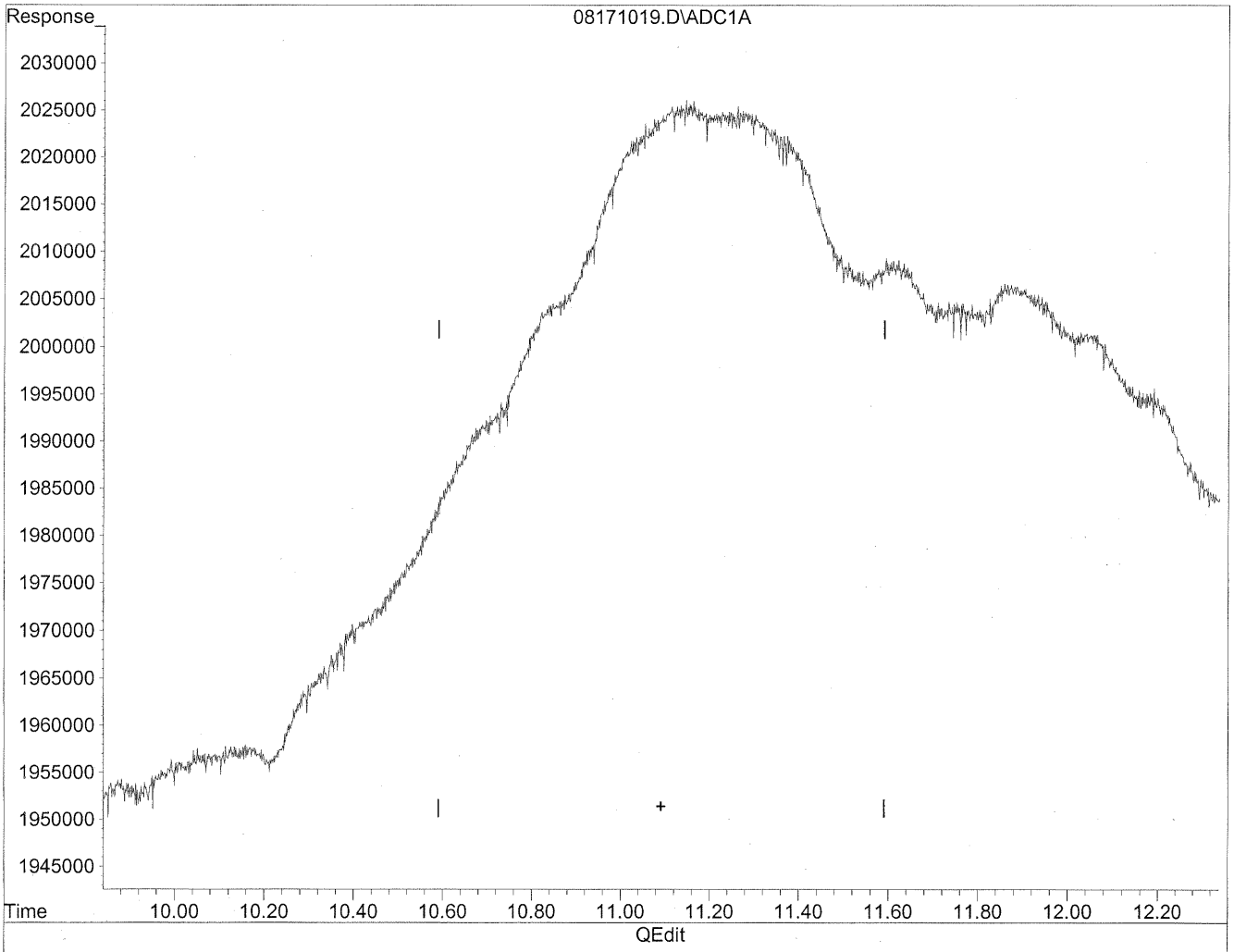
(+) = Expected Retention Time

08171019.D TO110709.M Sat Aug 22 14:37:53 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171019.D Vial: 32
Acq On : 18 Aug 2009 8:24 pm Operator: HC
Sample : P0902772-009 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(12) 2,5-Dimethylbenzaldehyde

0.00min 0.000ng/ml d

response 0

HC
8/22/09
not read

8/23/09

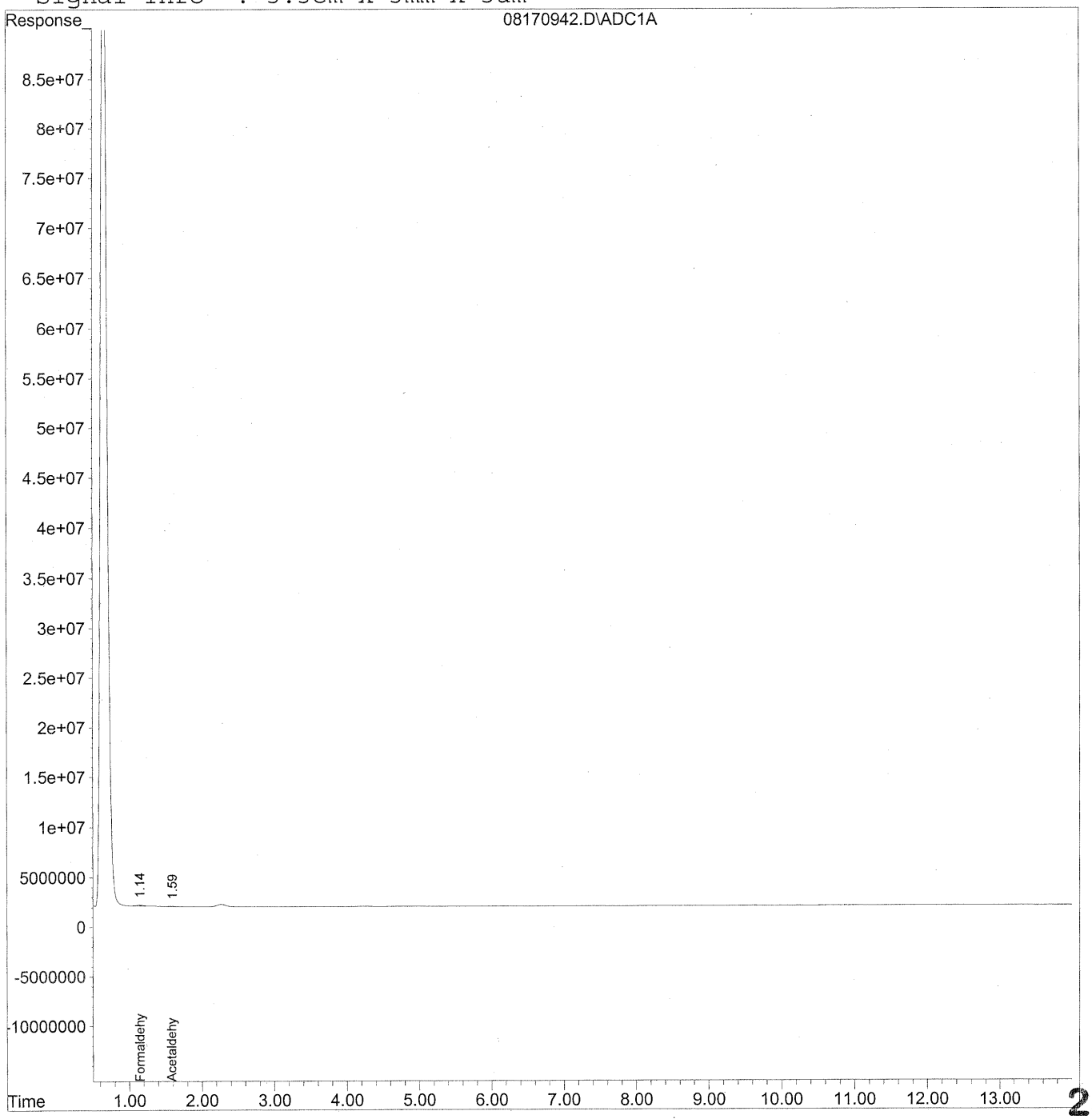
236

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_08\17\08170942.D Vial: 41
Acq On : 18 Aug 2009 1:07 am Operator: HC
Sample : P0902772-009 back1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:21 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



237

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170942.D Vial: 41
Acq On : 18 Aug 2009 1:07 am Operator: HC
Sample : P0902772-009 back1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:21 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

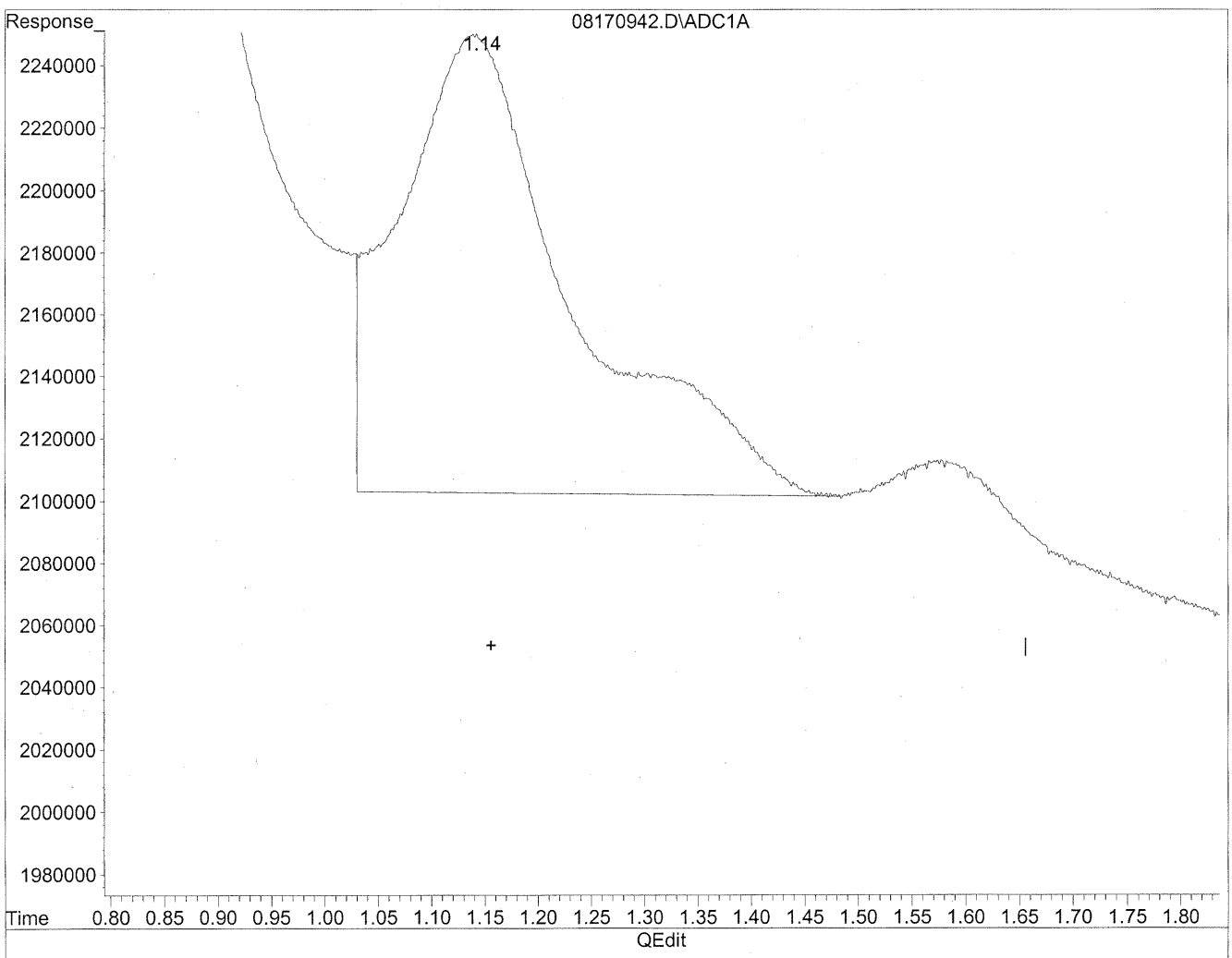
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.14	5518603	30.061	ng/mlm
2)	Acetaldehyde	1.59	1924004	13.721	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170942.D Vial: 41
Acq On : 18 Aug 2009 1:07 am Operator: HC
Sample : P0902772-009 back1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.14min 91.106ng/ml

response 16725319

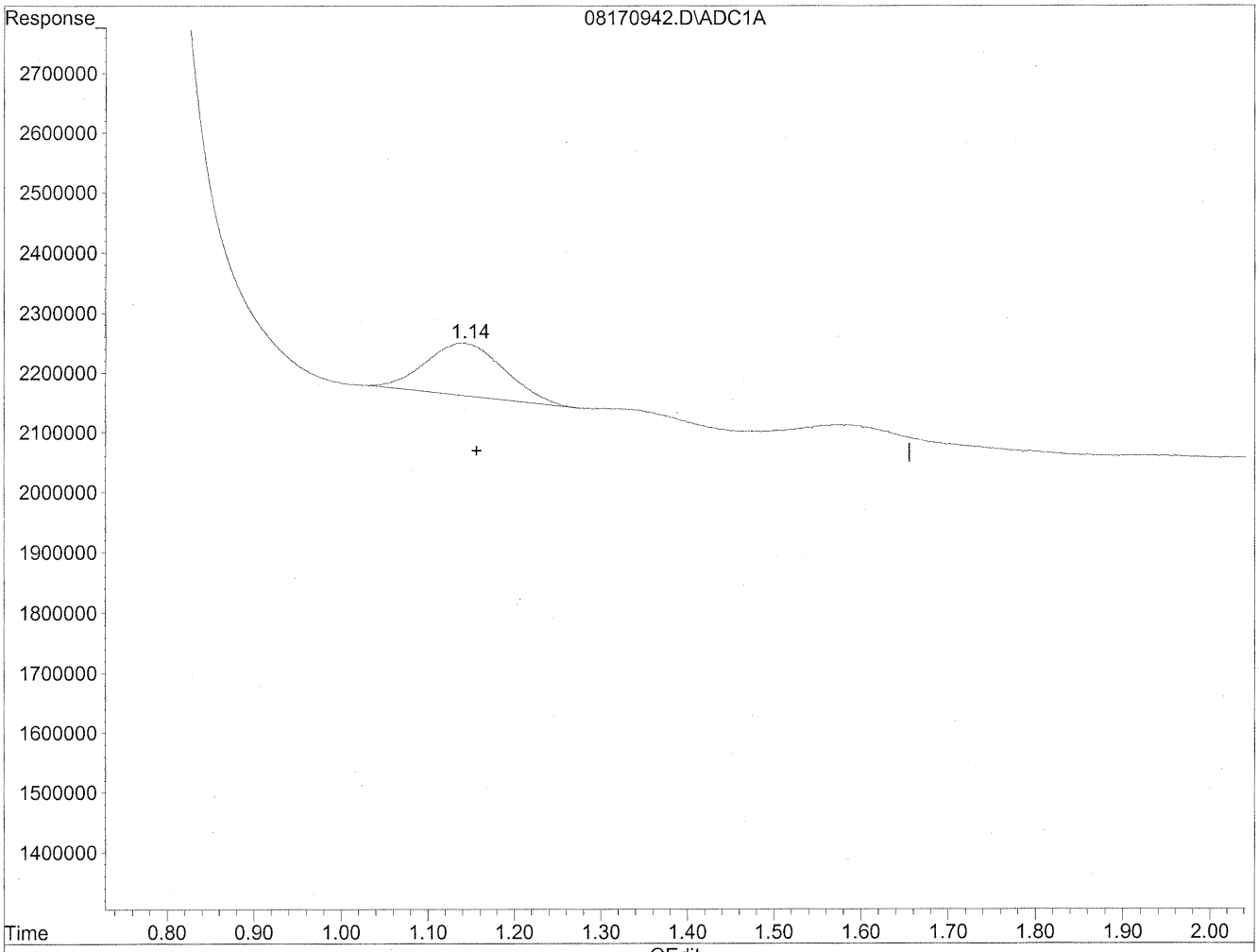
(+) = Expected Retention Time

08170942.D TO110709.M Fri Aug 21 18:20:44 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170942.D Vial: 41
Acq On : 18 Aug 2009 1:07 am Operator: HC
Sample : P0902772-009 back1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.14min 30.061ng/ml m

response 5518603

HC 8/24/09 LC

KB 8/25/09

240

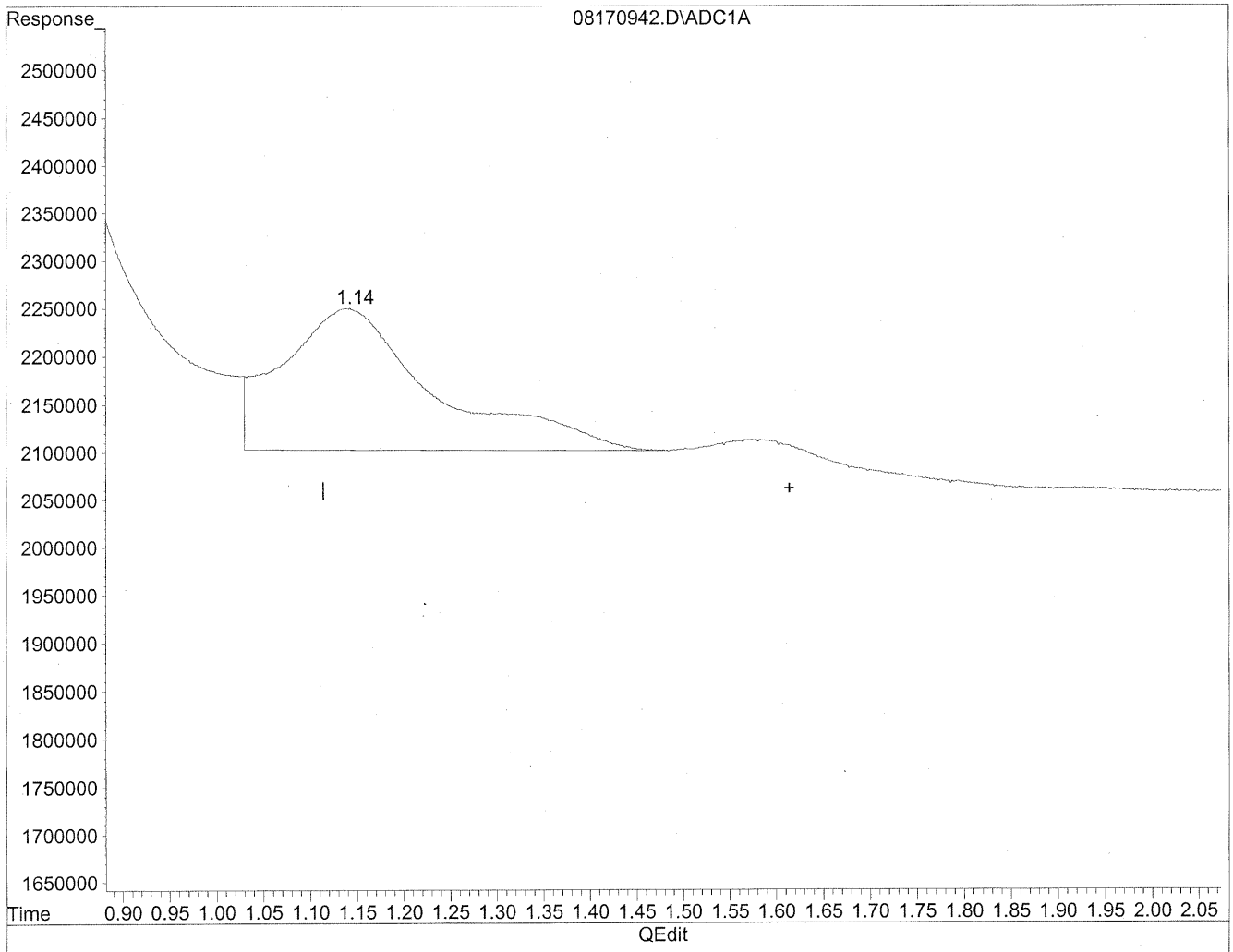
(+) = Expected Retention Time

08170942.D TO110709.M Fri Aug 21 18:20:53 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170942.D Vial: 41
Acq On : 18 Aug 2009 1:07 am Operator: HC
Sample : P0902772-009 back1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.14min 119.276ng/ml

response 16725319

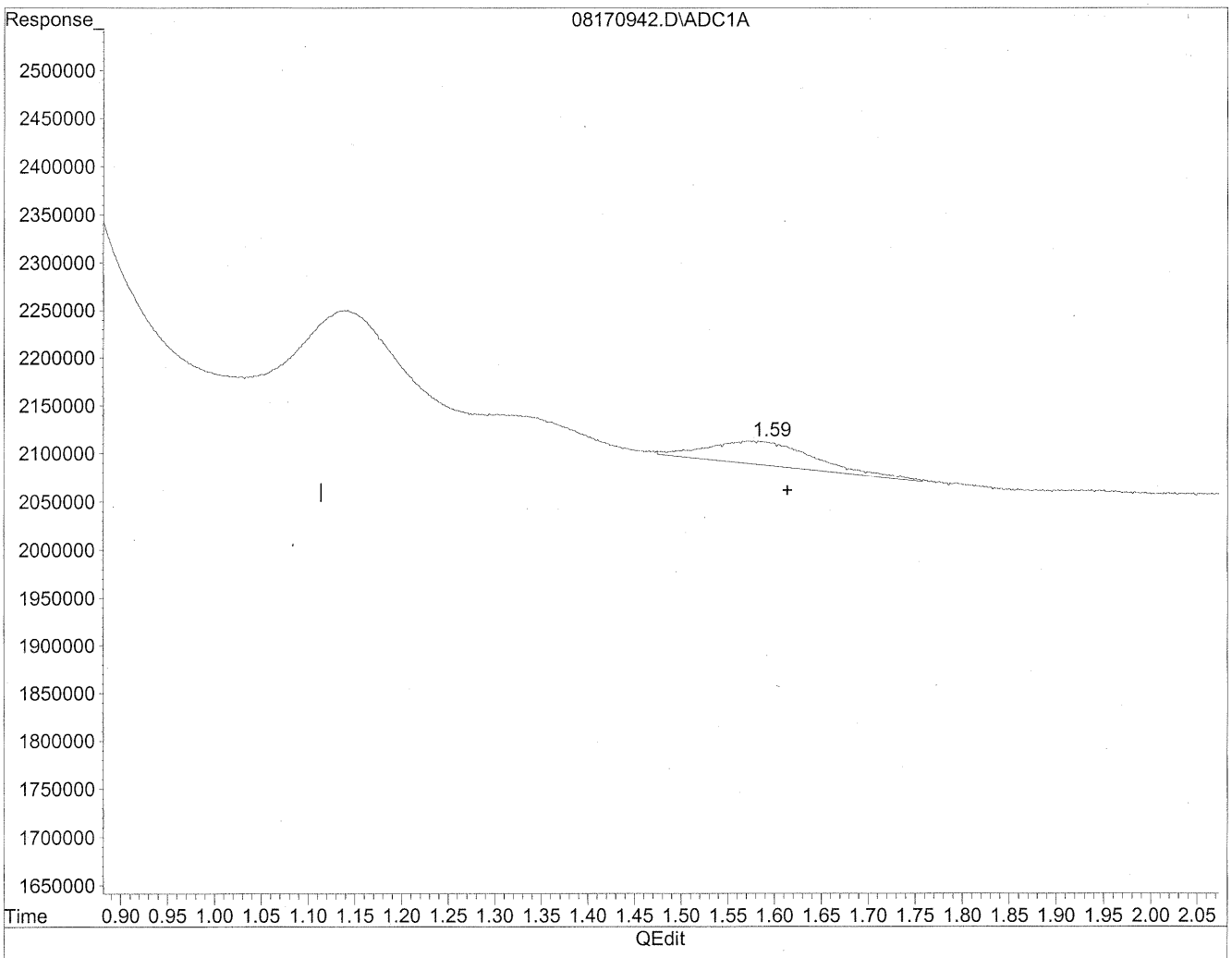
(+) = Expected Retention Time

08170942.D TO110709.M Fri Aug 21 18:21:01 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170942.D Vial: 41
Acq On : 18 Aug 2009 1:07 am Operator: HC
Sample : P0902772-009 back1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.59min 13.721ng/ml m

response 1924004

HC 8/21/09 WP

KR 8/23/09

242

(+) = Expected Retention Time

08170942.D TO110709.M Fri Aug 21 18:21:13 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **100451**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P0902772-010

Test Code: EPA Method TO-11A
Instrument ID: Waters LC Module I Plus/UV_Vis 360/LC1
Analyst: Hani Cherazaie
Sampling Media: Silica Gel DNPH Tube
Test Notes: **BC**

Date Collected: 8/10/09
Date Received: 8/12/09
Date Analyzed: 8/18/09
Desorption Volume: 1.0 ml
Volume Sampled: 110.9 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**410**	**3.7**	0.90	**3.0**	0.73	
75-07-0	**Acetaldehyde**	**190**	**1.7**	0.90	**0.96**	0.50	
123-38-6	Propionaldehyde	< 100	ND	0.90	ND	0.38	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.90	ND	0.31	
123-72-8	Butyraldehyde	< 100	ND	0.90	ND	0.31	
100-52-7	Benzaldehyde	< 100	ND	0.90	ND	0.21	
590-86-3	Isovaleraldehyde	< 100	ND	0.90	ND	0.26	
110-62-3	Valeraldehyde	< 100	ND	0.90	ND	0.26	
529-20-4	o-Tolualdehyde	< 100	ND	0.90	ND	0.18	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	1.8	ND	0.37	
66-25-1	n-Hexaldehyde	< 100	ND	0.90	ND	0.22	
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	ND	0.90	ND	0.16	

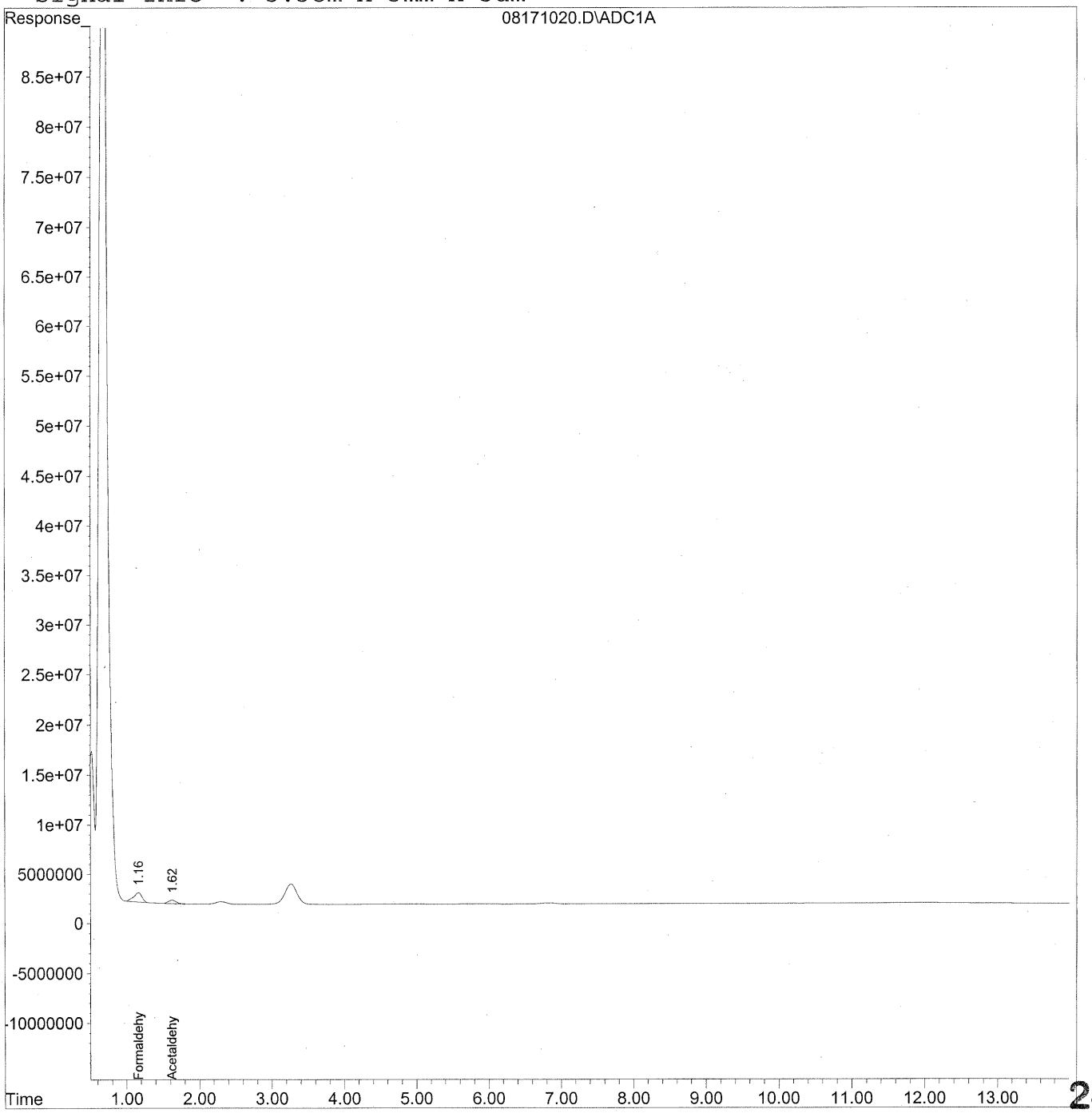
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.

Verified By: _____ Date: 8/25/09

Data File : J:\LC01\DATA\TO11\2009_08\17\08171020.D Vial: 33
Acq On : 18 Aug 2009 8:39 pm Operator: HC
Sample : P0902772-010 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:39 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



244

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171020.D Vial: 33
Acq On : 18 Aug 2009 8:39 pm Operator: HC
Sample : P0902772-010 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:39 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

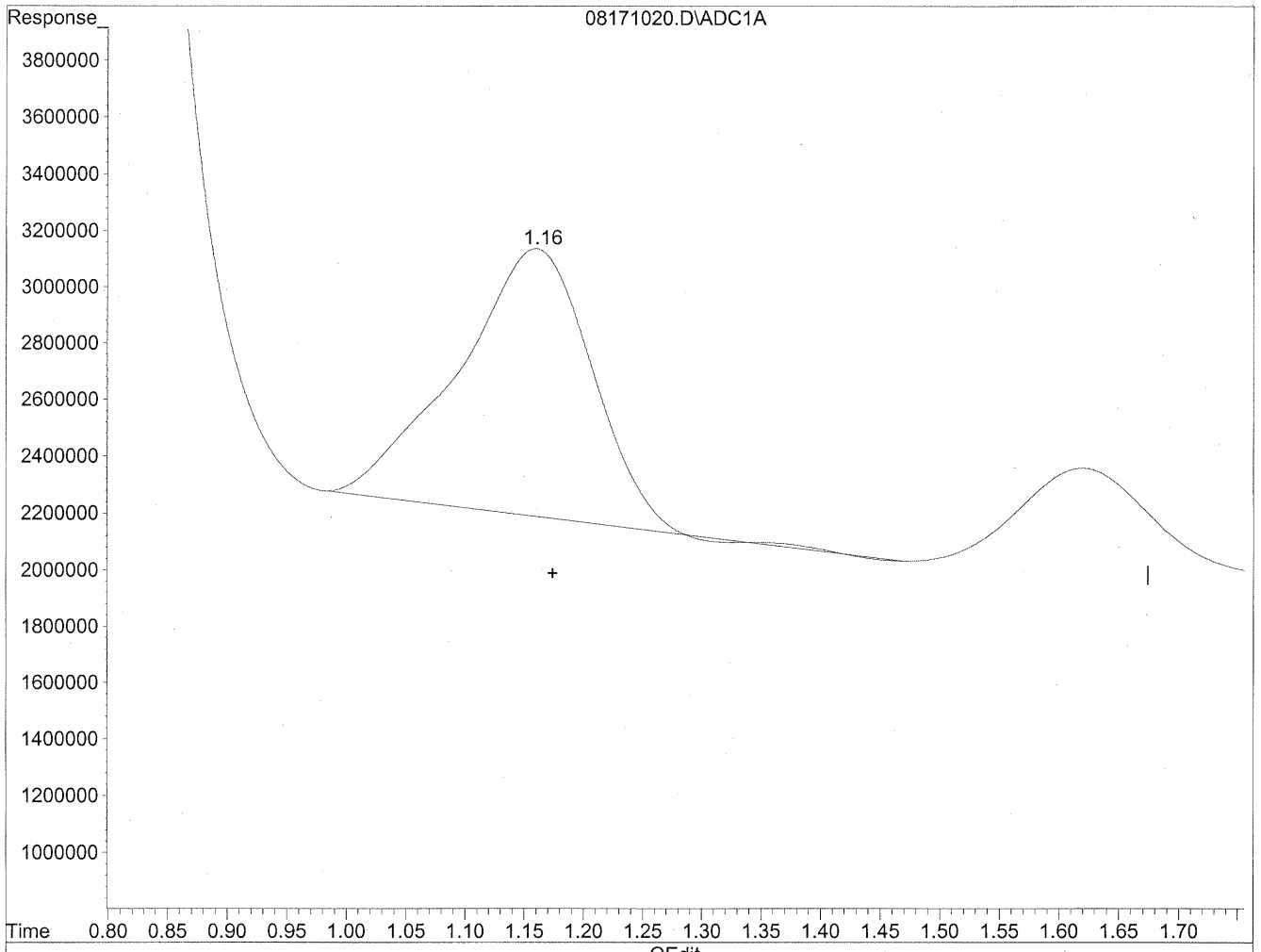
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	75269493	410.006	ng/mlm
2)	Acetaldehyde	1.62	26816472	191.241	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/mld
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171020.D Vial: 33
Acq On : 18 Aug 2009 8:39 pm Operator: HC
Sample : P0902772-010 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.16min 403.715ng/ml

response 74114523

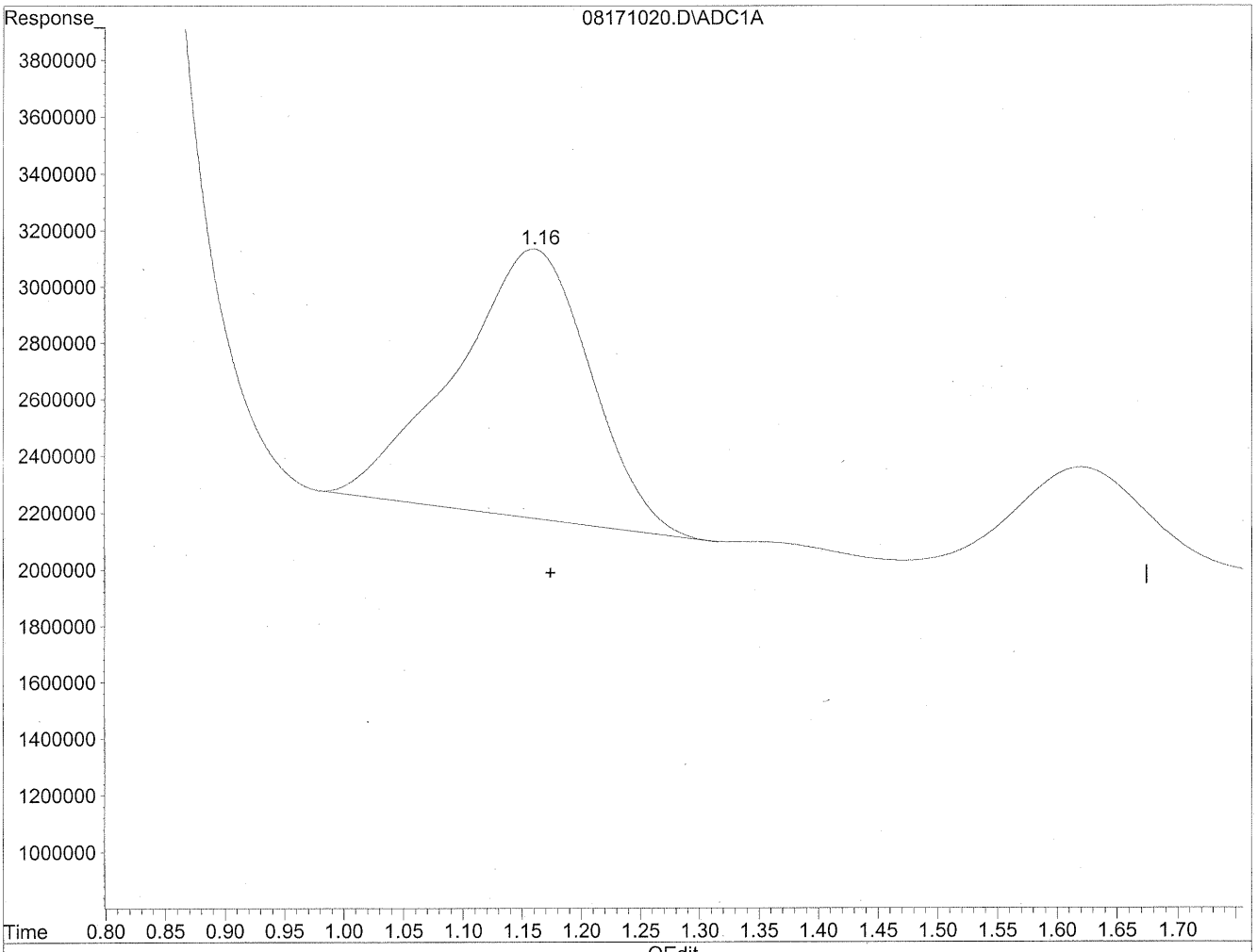
246

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171020.D Vial: 33
Acq On : 18 Aug 2009 8:39 pm Operator: HC
Sample : P0902772-010 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.16min 410.006ng/ml m

response 75269493

HC 8/22/09
LC
KP 8/23/09

247

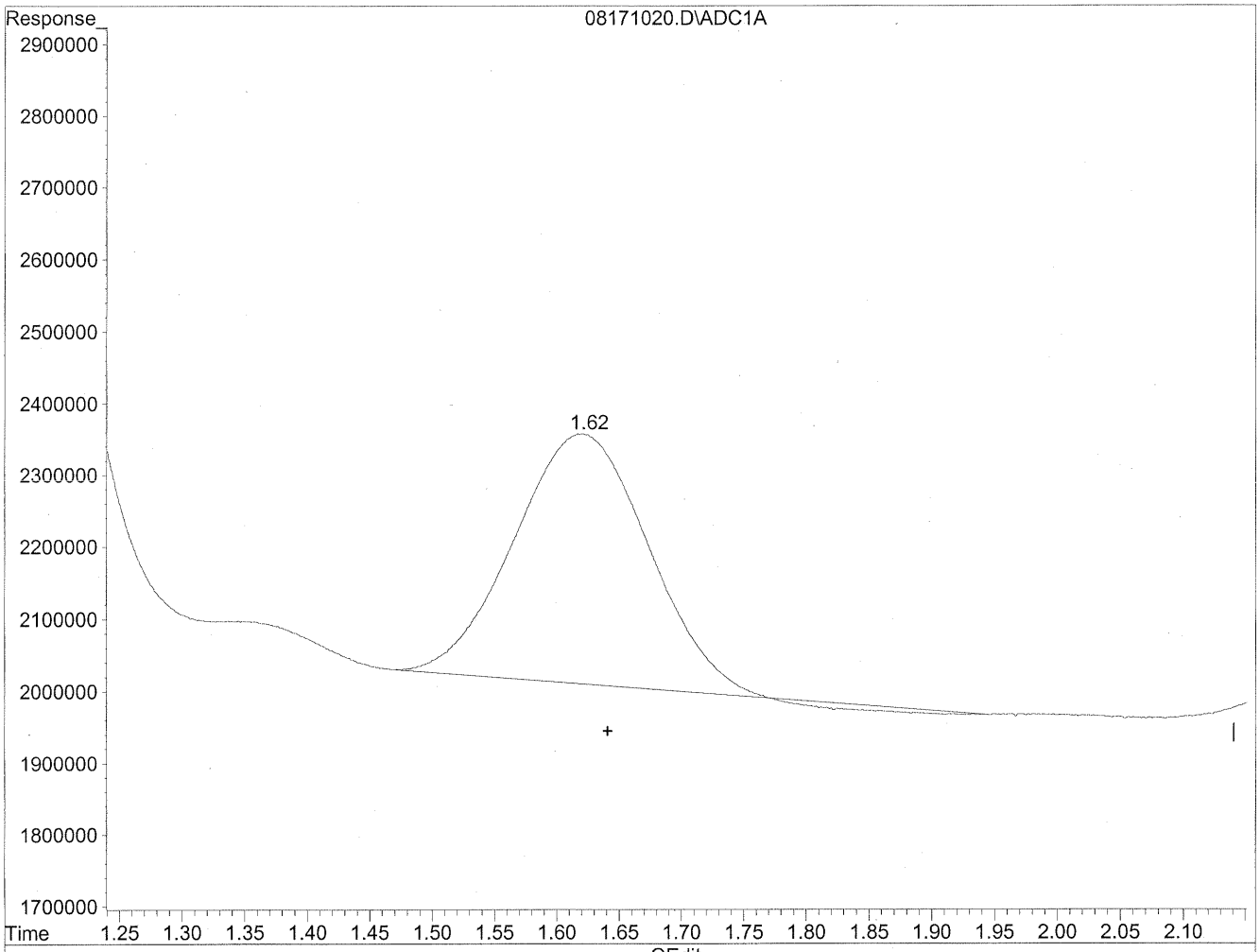
(+) = Expected Retention Time

08171020.D TO110709.M Sat Aug 22 14:38:34 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171020.D Vial: 33
Acq On : 18 Aug 2009 8:39 pm Operator: HC
Sample : P0902772-010 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.62min 181.830ng/ml

response 25496889

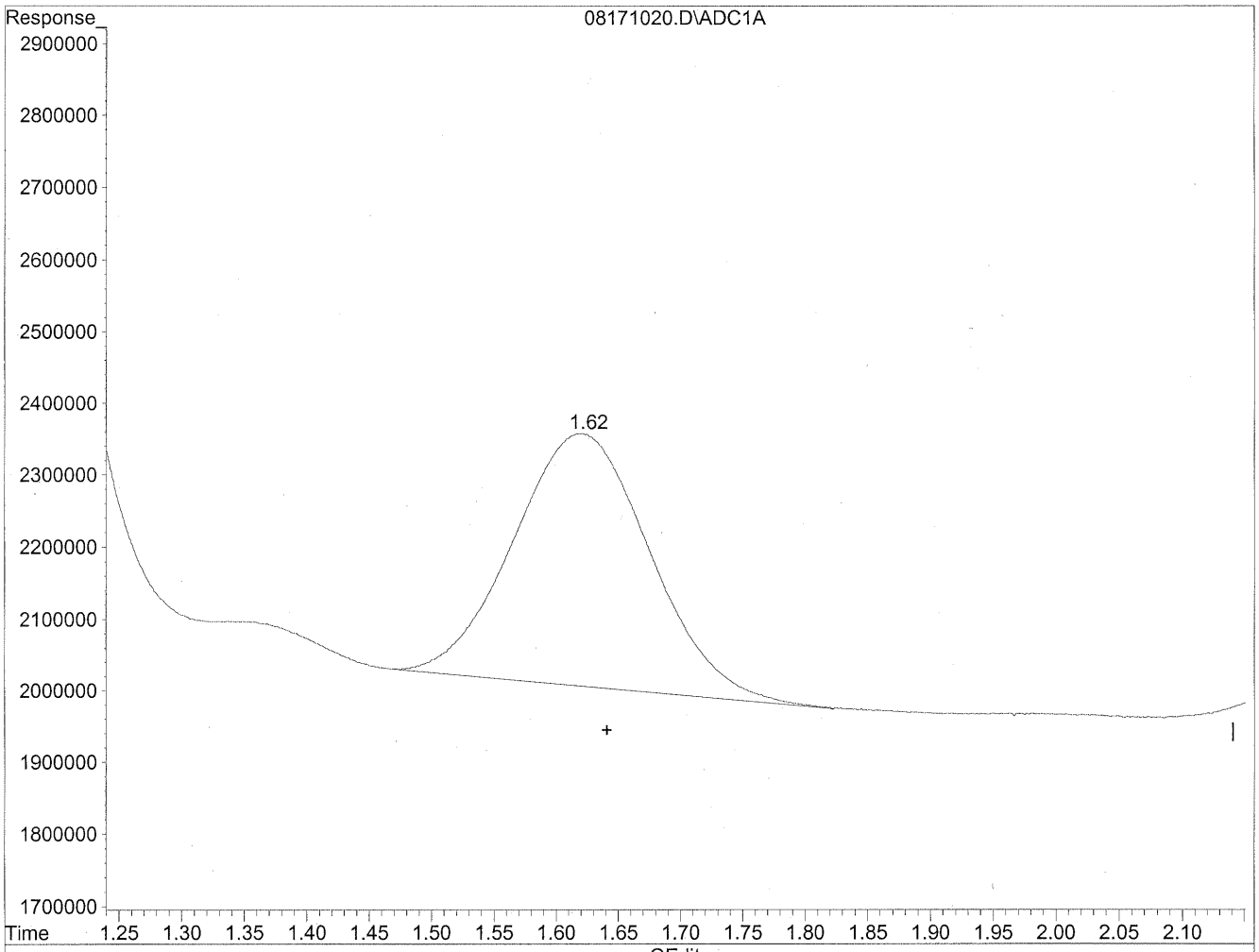
(+) = Expected Retention Time

08171020.D TO110709.M Sat Aug 22 14:39:02 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171020.D Vial: 33
Acq On : 18 Aug 2009 8:39 pm Operator: HC
Sample : P0902772-010 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.62min 191.241ng/ml m

response 26816472

HC
8/22/09
IC

KP 8/23/09

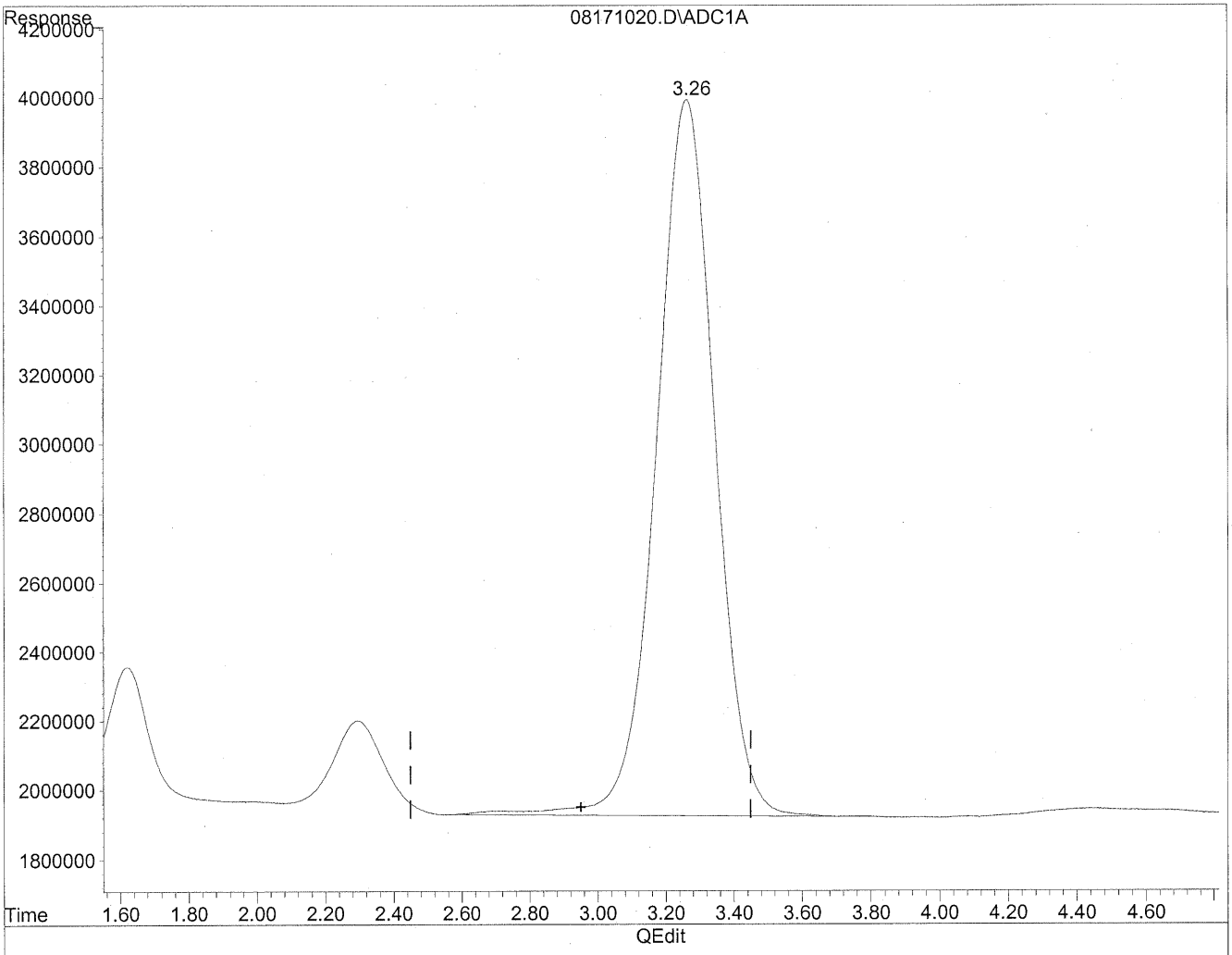
249

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171020.D Vial: 33
Acq On : 18 Aug 2009 8:39 pm Operator: HC
Sample : P0902772-010 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.26min 2320.301ng/ml

response 247565064

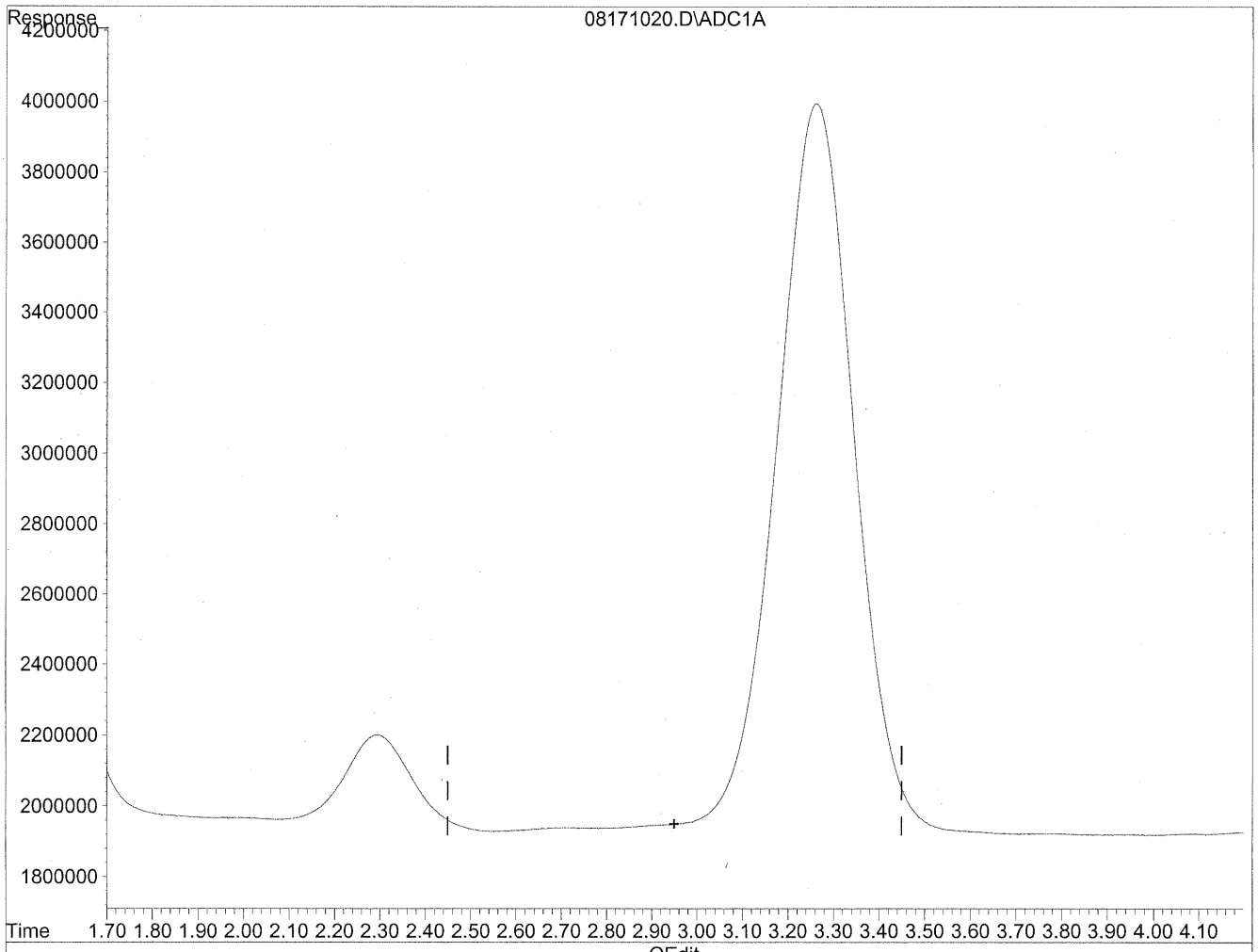
(+) = Expected Retention Time

08171020.D TO110709.M Sat Aug 22 14:39:11 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171020.D Vial: 33
Acq On : 18 Aug 2009 8:39 pm Operator: HC
Sample : P0902772-010 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09 MP

KP 8/23/09

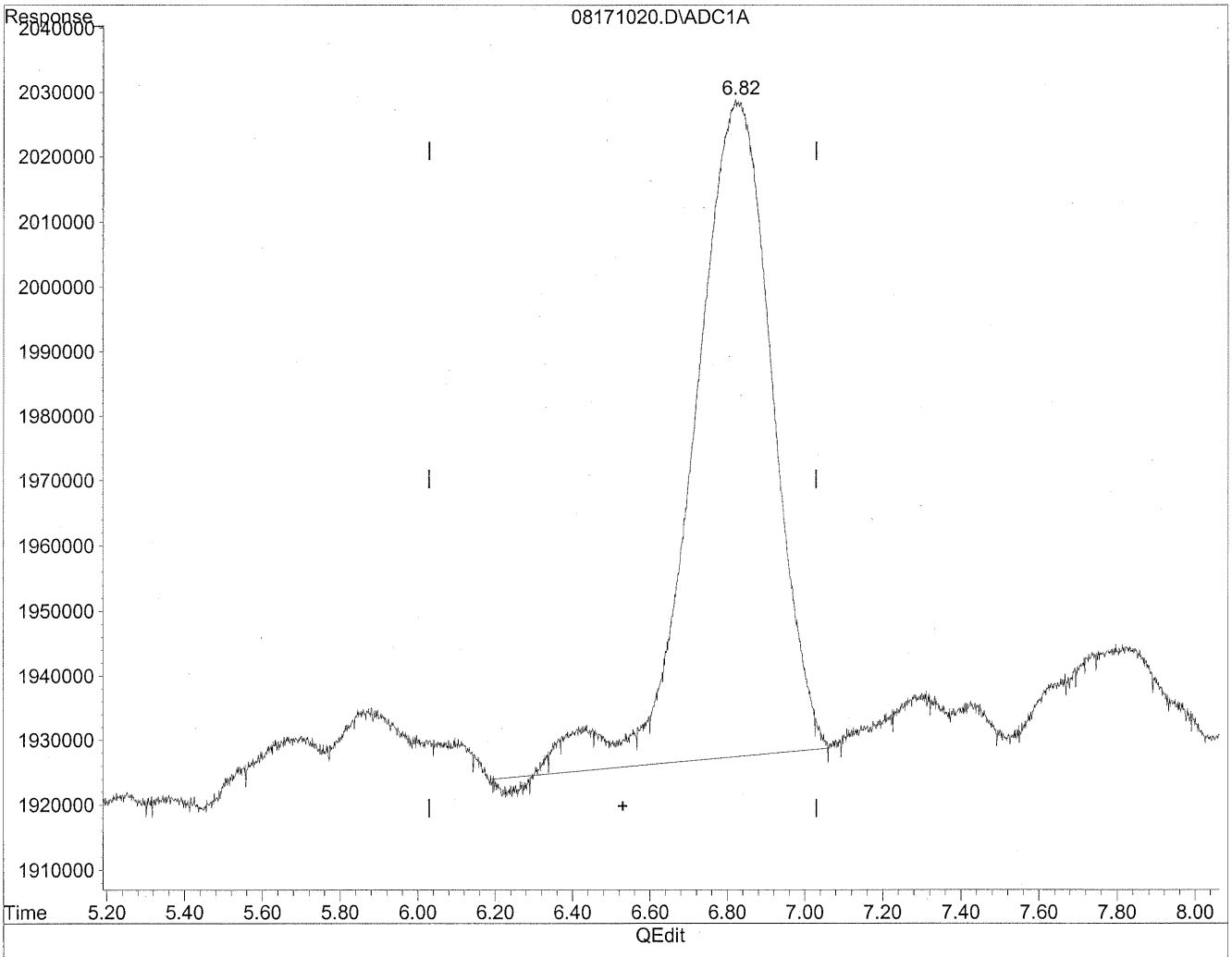
(+) = Expected Retention Time

08171020.D TO110709.M Sat Aug 22 14:39:13 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171020.D Vial: 33
Acq On : 18 Aug 2009 8:39 pm Operator: HC
Sample : P0902772-010 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.83min 217.551ng/ml

response 14329917

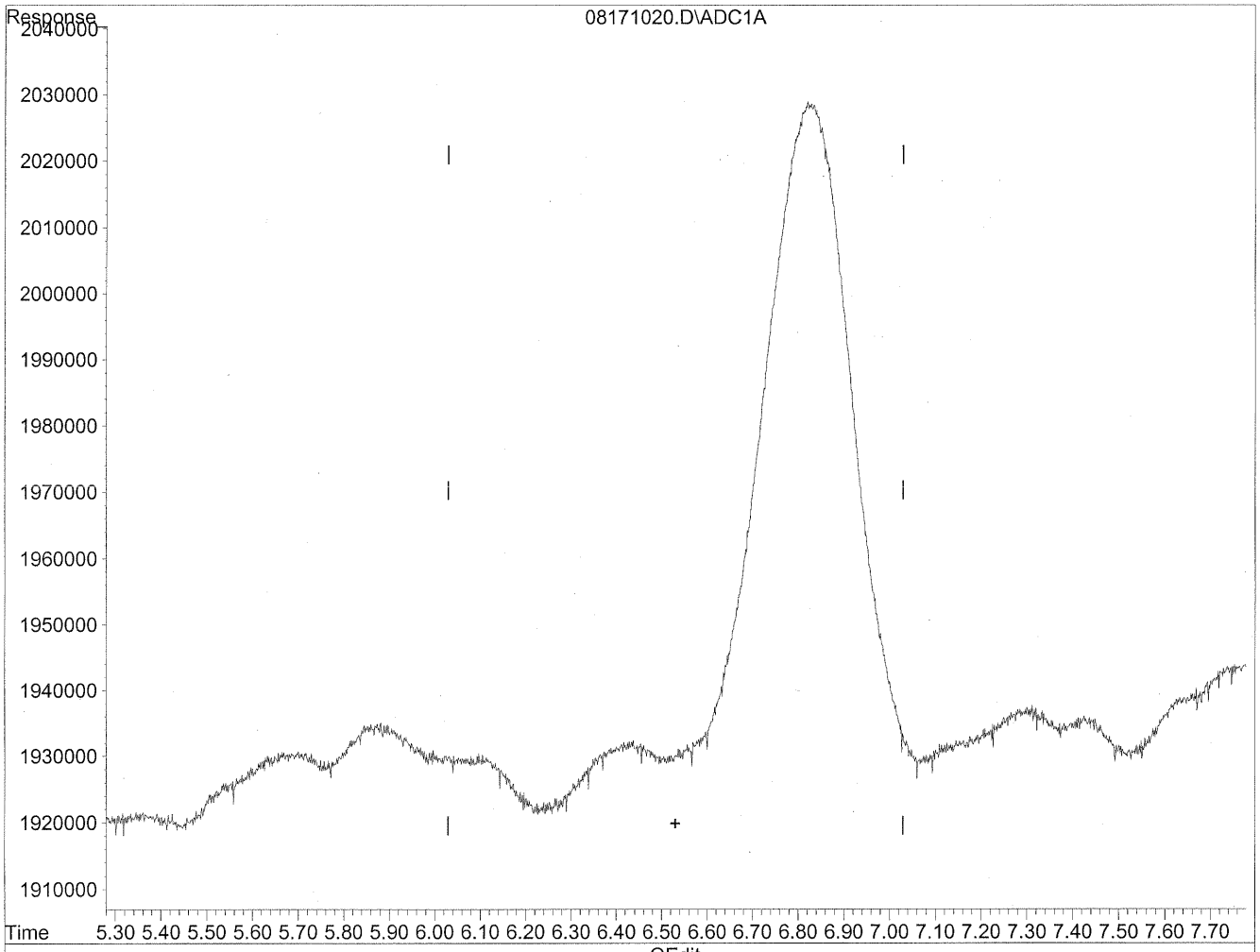
(+) = Expected Retention Time

08171020.D TO110709.M Sat Aug 22 14:39:22 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171020.D Vial: 33
Acq On : 18 Aug 2009 8:39 pm Operator: HC
Sample : P0902772-010 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(6) Benzaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09 MP

KE 8/23/09

253

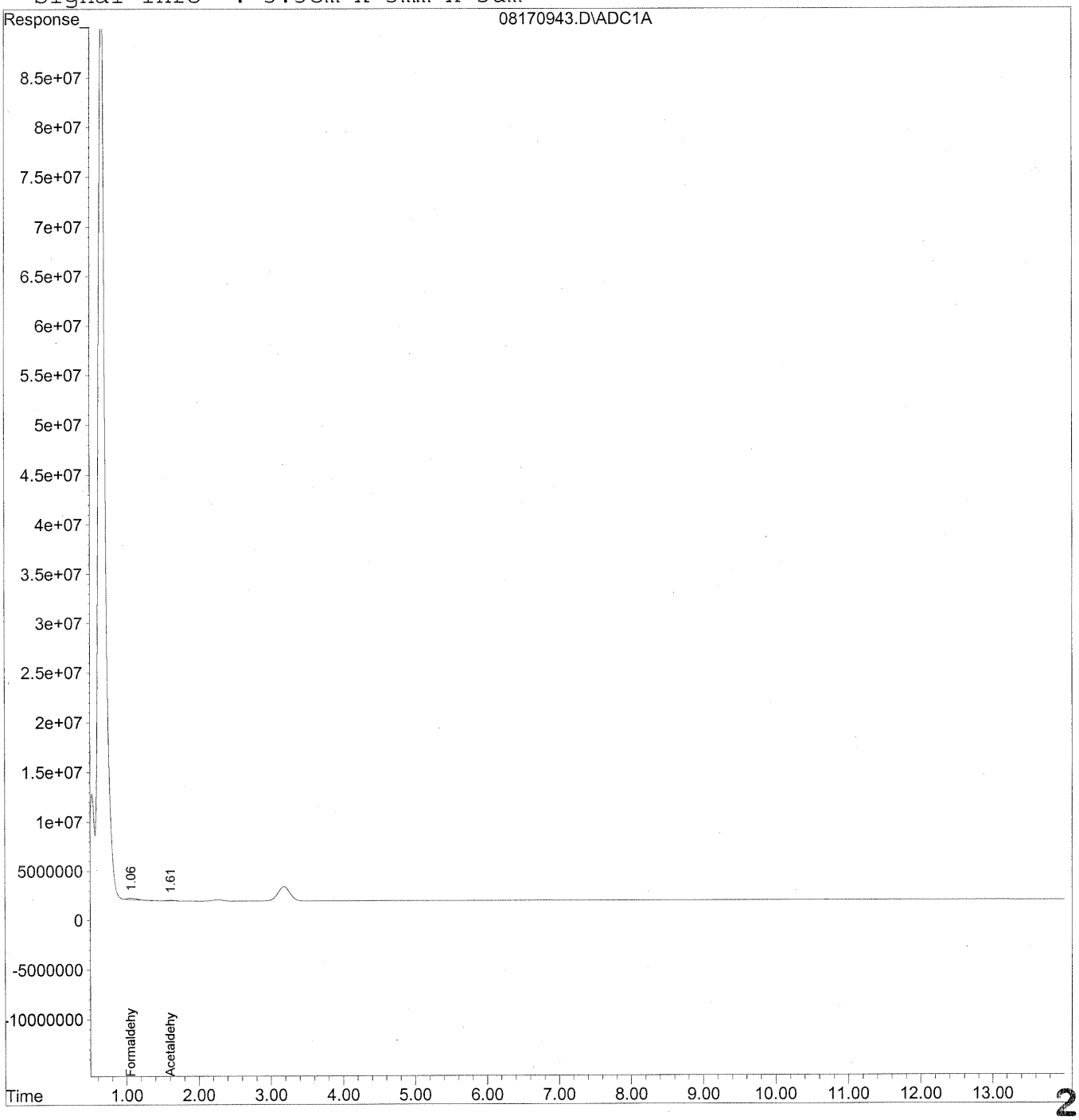
(+) = Expected Retention Time

08171020.D TO110709.M Sat Aug 22 14:39:25 2009

Data File : J:\LC01\DATA\TO11\2009_08\17\08170943.D Vial: 42
Acq On : 18 Aug 2009 1:22 am Operator: HC
Sample : P0902772-010 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:34 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



254

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170943.D Vial: 42
Acq On : 18 Aug 2009 1:22 am Operator: HC
Sample : P0902772-010 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:34 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

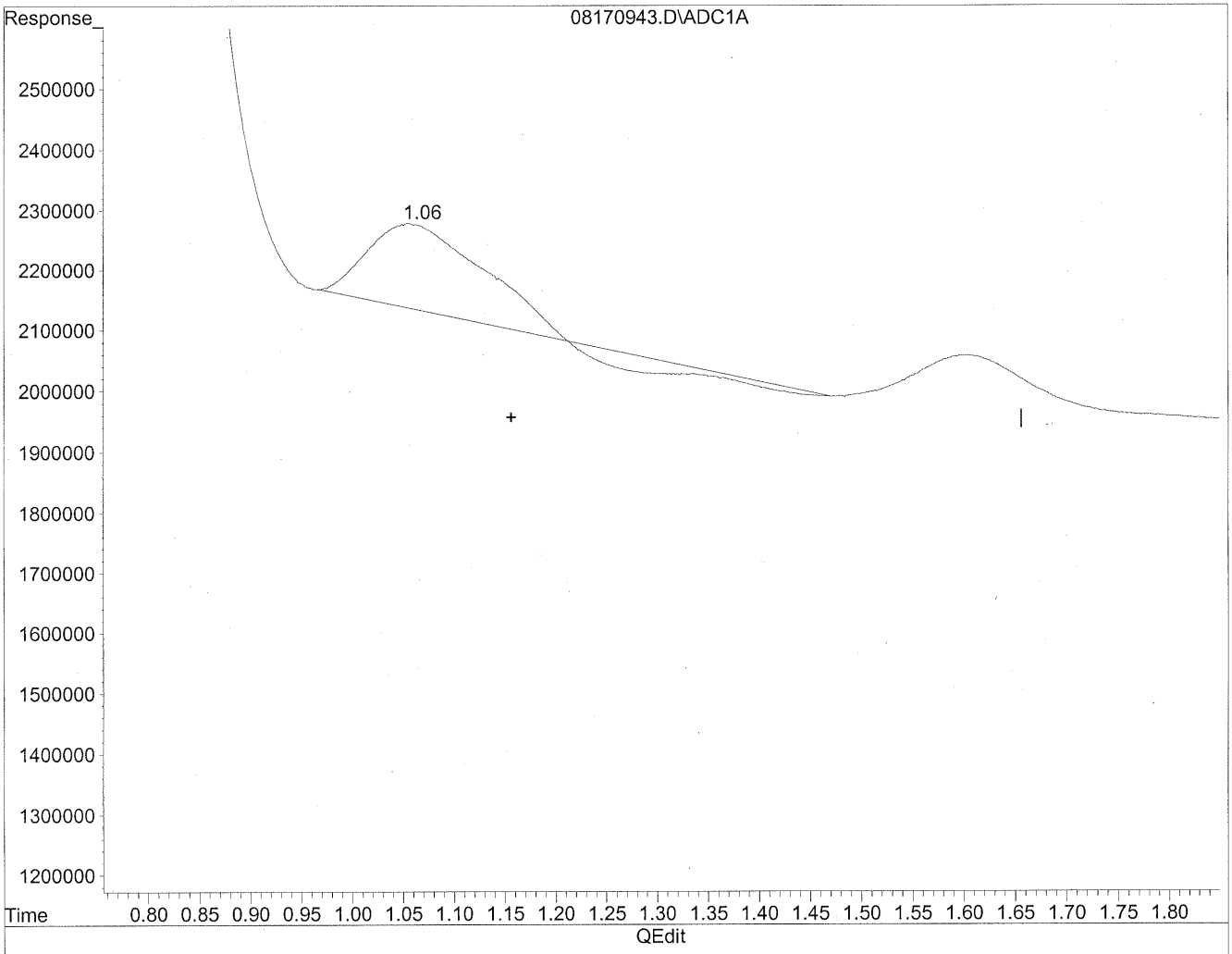
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.06	9550286	52.022	ng/ml
2)	Acetaldehyde	1.60	5493432	39.176	ng/ml
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170943.D Vial: 42
Acq On : 18 Aug 2009 1:22 am Operator: HC
Sample : P0902772-010 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.06min 52.022ng/ml

response 9550286

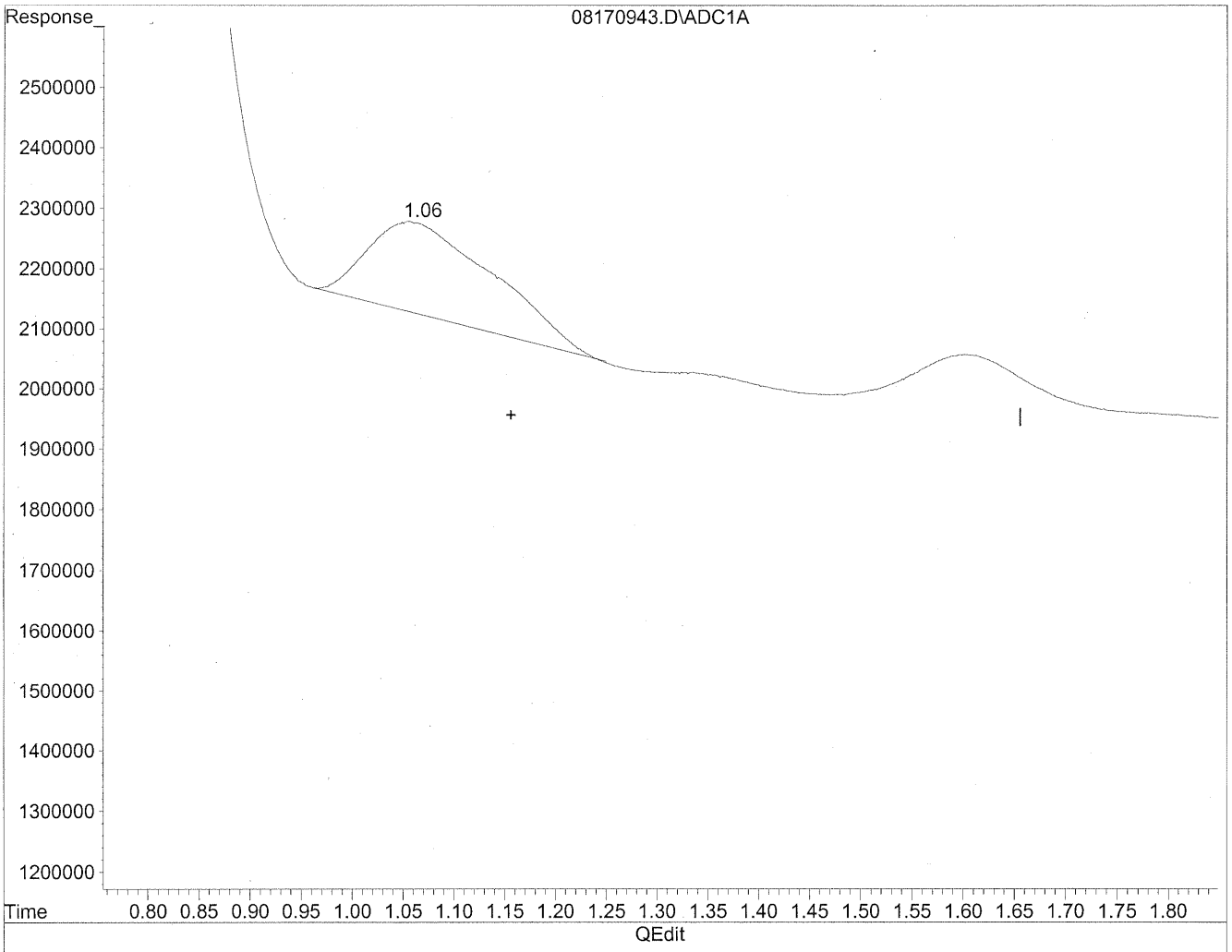
(+) = Expected Retention Time

08170943.D TO110709.M Fri Aug 21 18:21:31 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170943.D Vial: 42
Acq On : 18 Aug 2009 1:22 am Operator: HC
Sample : P0902772-010 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.06min 72.389ng/ml m

response 13289338

HC 8/21/09
TC

KE 8/23/09

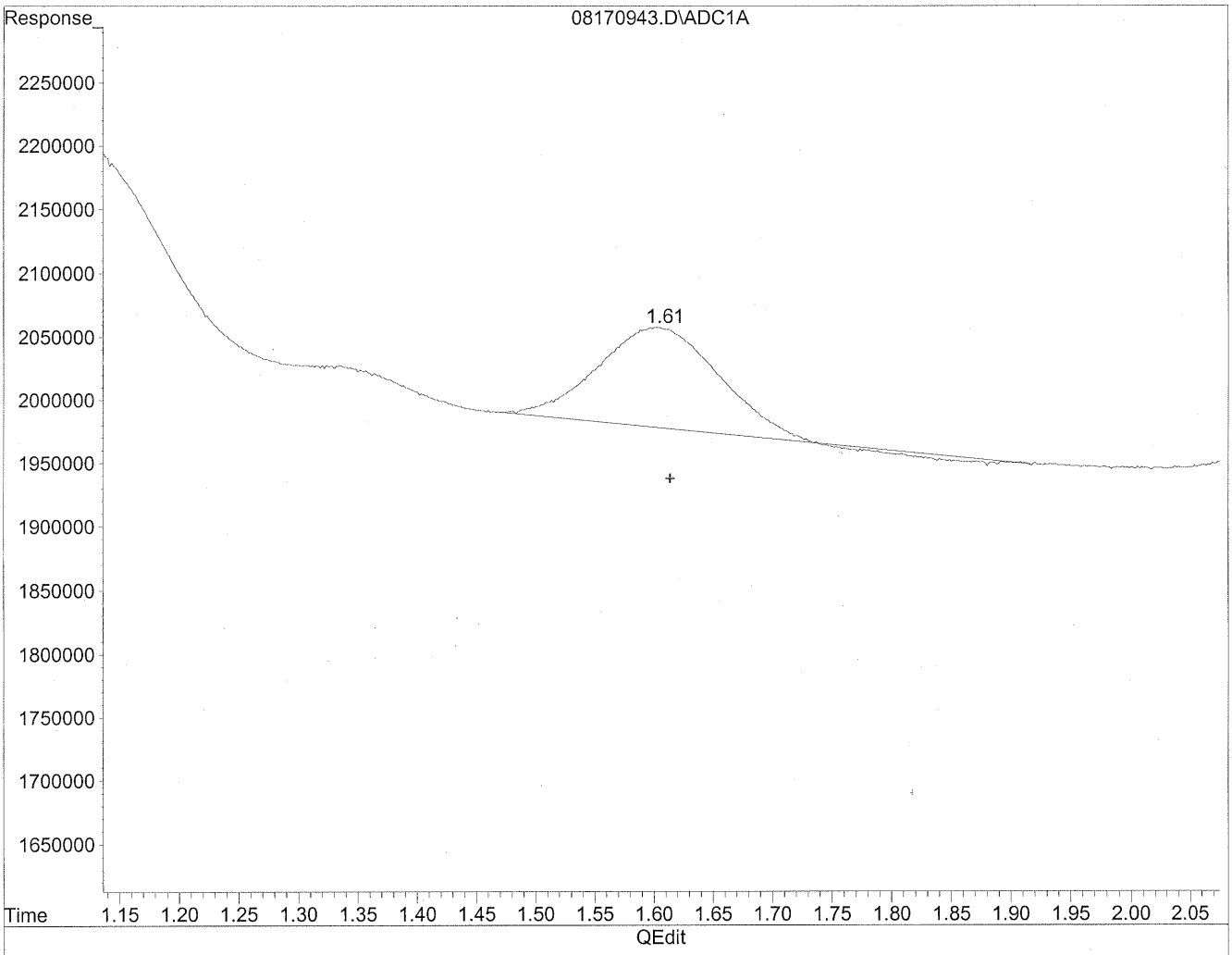
(+) = Expected Retention Time

08170943.D TO110709.M Fri Aug 21 18:21:38 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170943.D Vial: 42
Acq On : 18 Aug 2009 1:22 am Operator: HC
Sample : P0902772-010 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.60min 39.176ng/ml

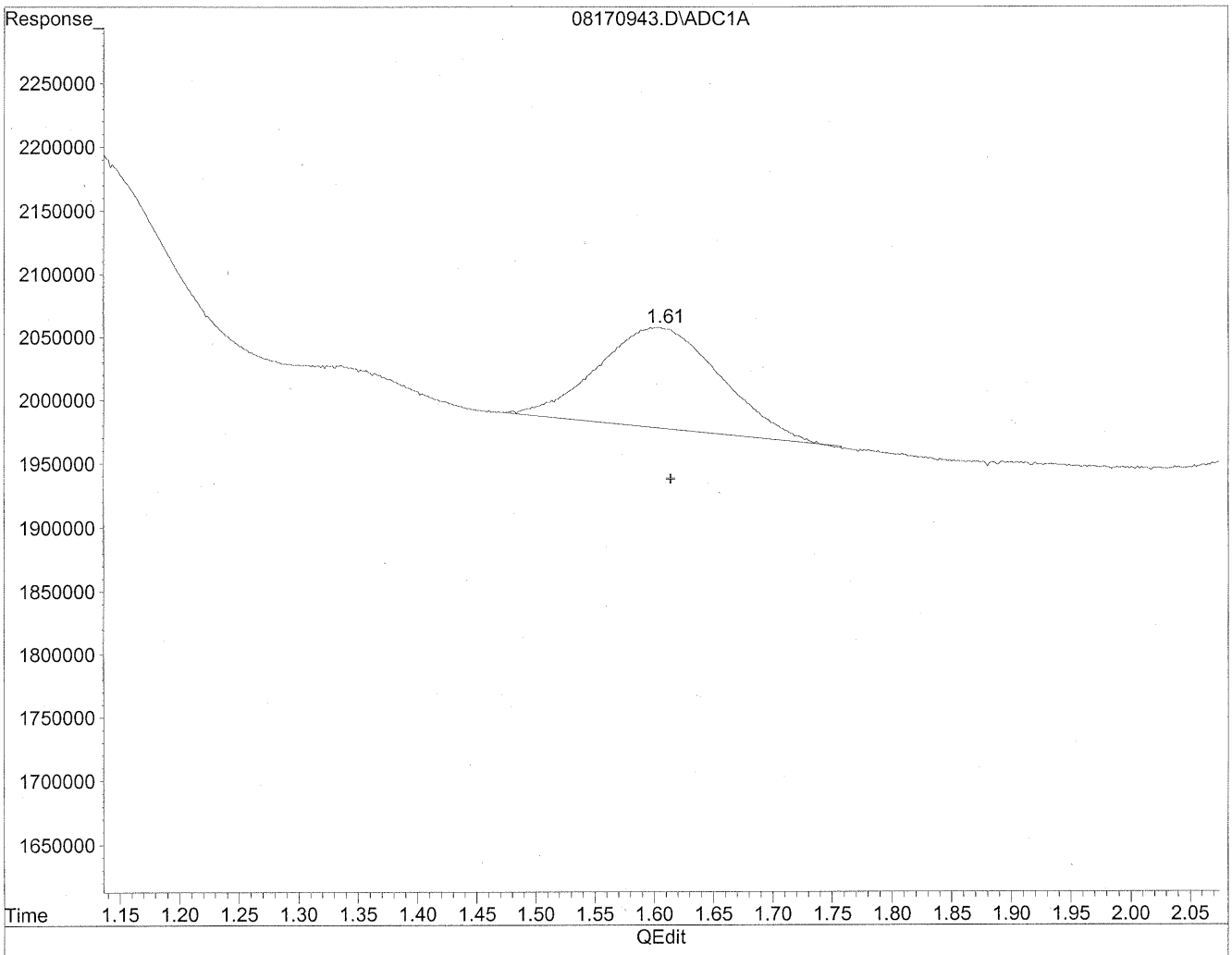
response 5493432

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170943.D Vial: 42
Acq On : 18 Aug 2009 1:22 am Operator: HC
Sample : P0902772-010 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:21 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.61min 40.966ng/ml m

response 5744450

HC 8/24/09 LC

KB 8/23/09

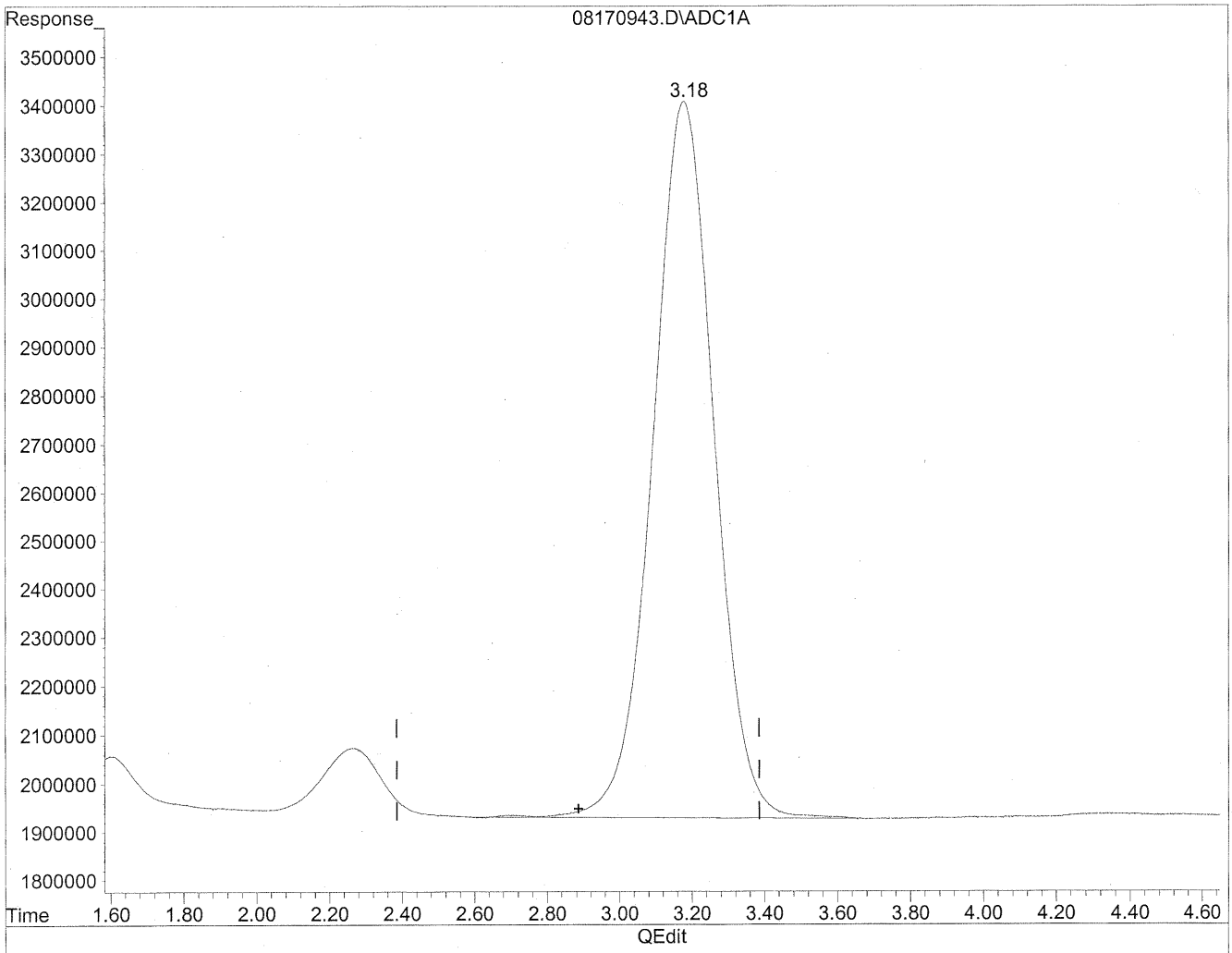
(+) = Expected Retention Time

08170943.D TO110709.M Fri Aug 21 18:22:03 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170943.D Vial: 42
Acq On : 18 Aug 2009 1:22 am Operator: HC
Sample : P0902772-010 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:33 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.18min 1625.770ng/ml

response 173461899

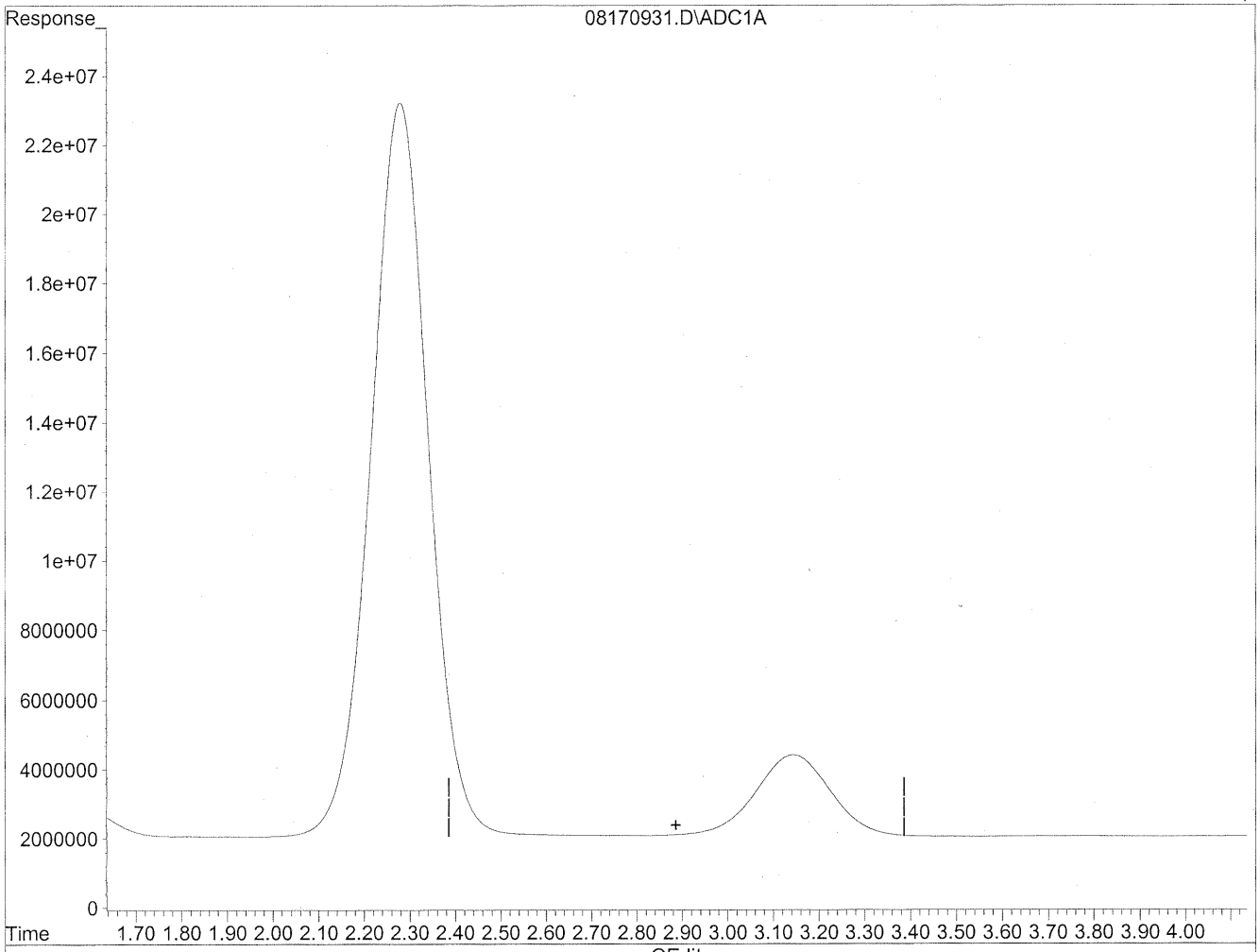
(+) = Expected Retention Time

08170943.D TO110709.M Sat Aug 22 12:34:43 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170931.D Vial: 30
Acq On : 17 Aug 2009 10:21 pm Operator: HC
Sample : P0902772-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:13 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/25/09 MP

261

(+) = Expected Retention Time

08170931.D TO110709.M Tue Aug 25 07:48:57 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **100457**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P0902772-011

Test Code: EPA Method TO-11A
Instrument ID: Waters LC Module I Plus/UV_Vis 360/LC1
Analyst: Hani Cherazaie
Sampling Media: Silica Gel DNPH Tube
Test Notes: **BC**

Date Collected: 8/10/09
Date Received: 8/12/09
Date Analyzed: 8/18 - 8/19/09
Desorption Volume: 1.0 ml
Volume Sampled: 100.6 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**11,000**	**110**	0.99	**91**	0.81	
75-07-0	**Acetaldehyde**	**14,000**	**140**	0.99	**76**	0.55	**BT**
123-38-6	**Propionaldehyde**	**790**	**7.8**	0.99	**3.3**	0.42	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.99	ND	0.35	
123-72-8	**Butyraldehyde**	**480**	**4.8**	0.99	**1.6**	0.34	
100-52-7	**Benzaldehyde**	**1,100**	**11**	0.99	**2.6**	0.23	
590-86-3	**Isovaleraldehyde**	**240**	**2.4**	0.99	**0.67**	0.28	
110-62-3	**Valeraldehyde**	**900**	**8.9**	0.99	**2.5**	0.28	
529-20-4	o-Tolualdehyde	< 100	ND	0.99	ND	0.20	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	2.0	ND	0.40	
66-25-1	**n-Hexaldehyde**	**2,700**	**26**	0.99	**6.5**	0.24	**M**
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	ND	0.99	ND	0.18	

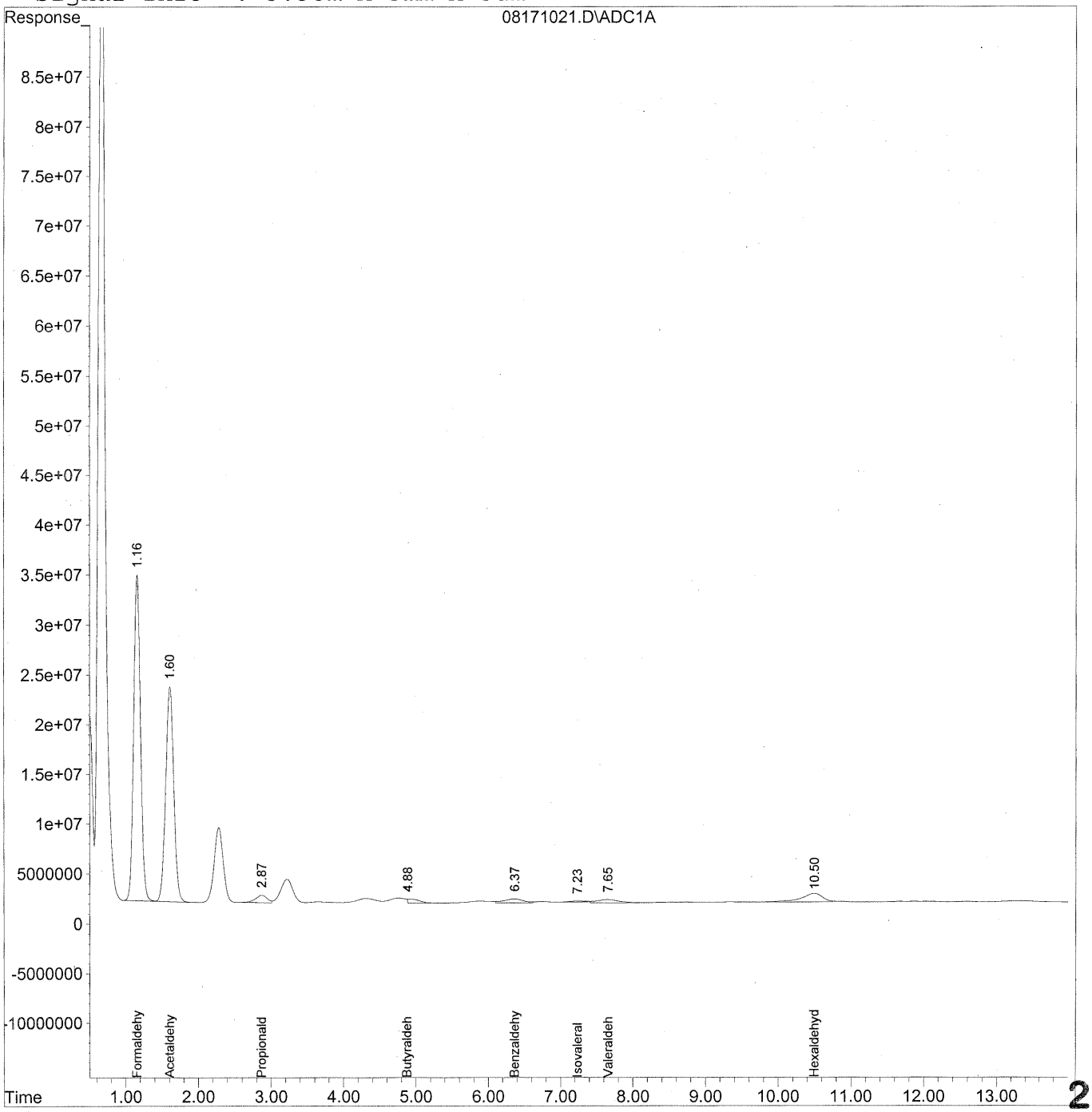
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
BT = Results indicated possible breakthrough; back section > 10% front section.
M = Matrix interference; results may be biased high.

Verified By: _____ Date: 8/25/09

Data File : J:\LC01\DATA\TO11\2009_08\17\08171021.D Vial: 34
Acq On : 18 Aug 2009 8:54 pm Operator: HC
Sample : P0902772-011 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:27 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



263

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171021.D Vial: 34
Acq On : 18 Aug 2009 8:54 pm Operator: HC
Sample : P0902772-011 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:27 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

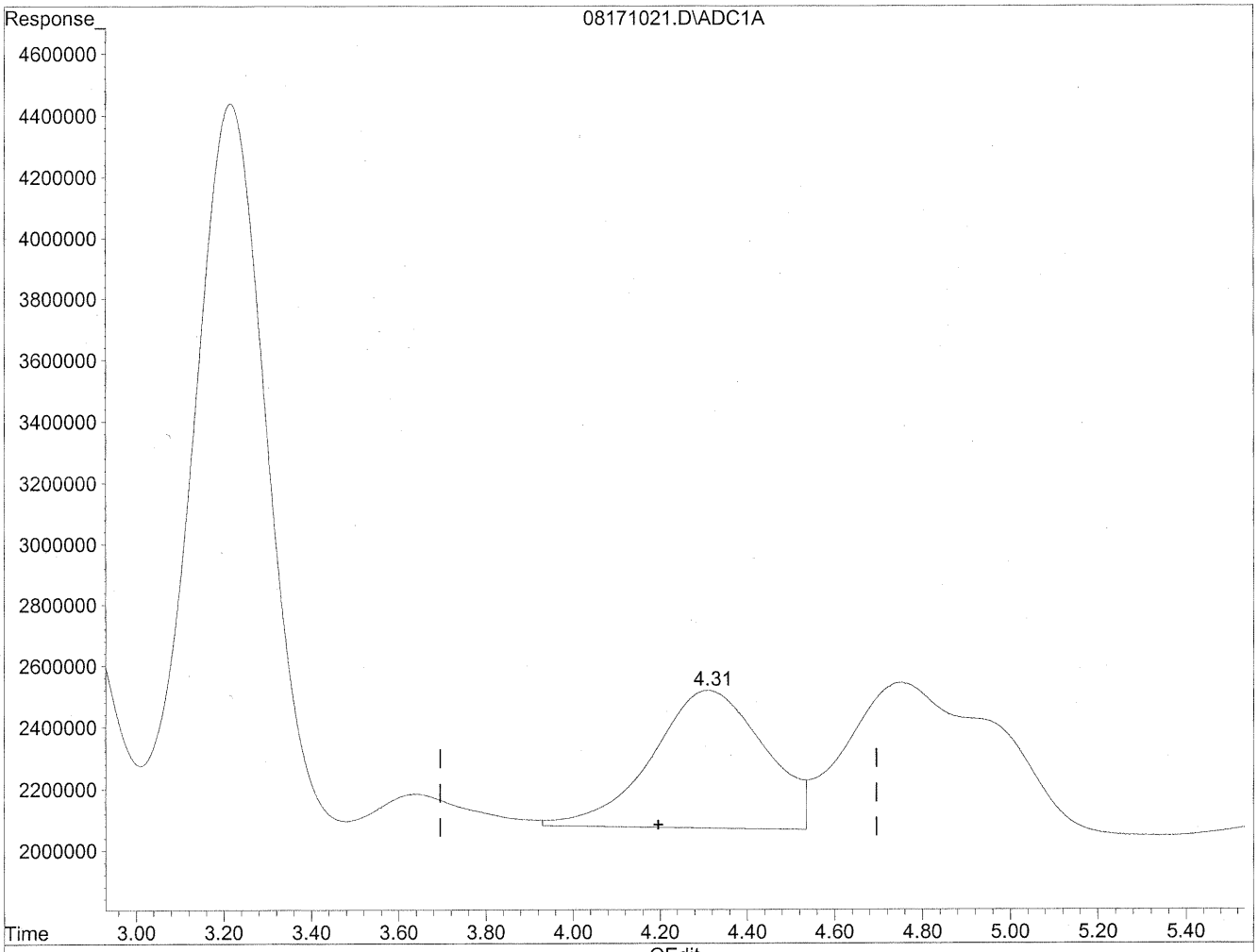
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	2124824580	11574.292	ng/ml
2)	Acetaldehyde	1.60	1694511676	12084.362	ng/ml
3)	Propionaldehyde	2.87	84166521	788.850	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	4.88	42238936	478.162	ng/mlm
6)	Benzaldehyde	6.37	73865582	1121.396	ng/mlm
7)	Isovaleraldehyde	7.23	18549852	237.056	ng/mlm
8)	Valeraldehyde	7.65	65949074	897.205	ng/mlm
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	10.50f	179026219	2658.393	ng/mlm
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171021.D Vial: 34
Acq On : 18 Aug 2009 8:54 pm Operator: HC
Sample : P0902772-011 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.31min 826.526ng/ml

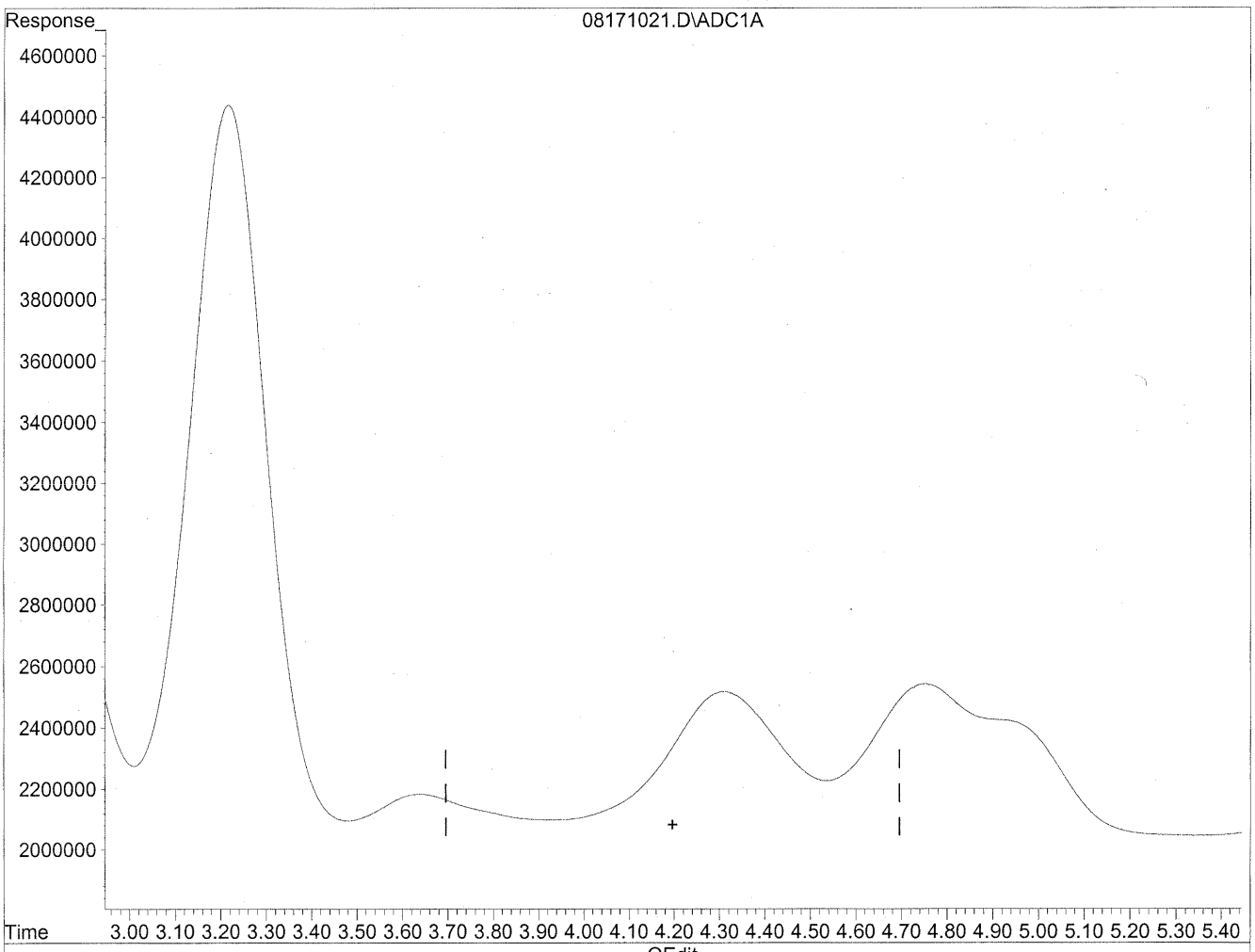
response 80516202

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171021.D Vial: 34
Acq On : 18 Aug 2009 8:54 pm Operator: HC
Sample : P0902772-011 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09 MP

HP 8/23/09

266

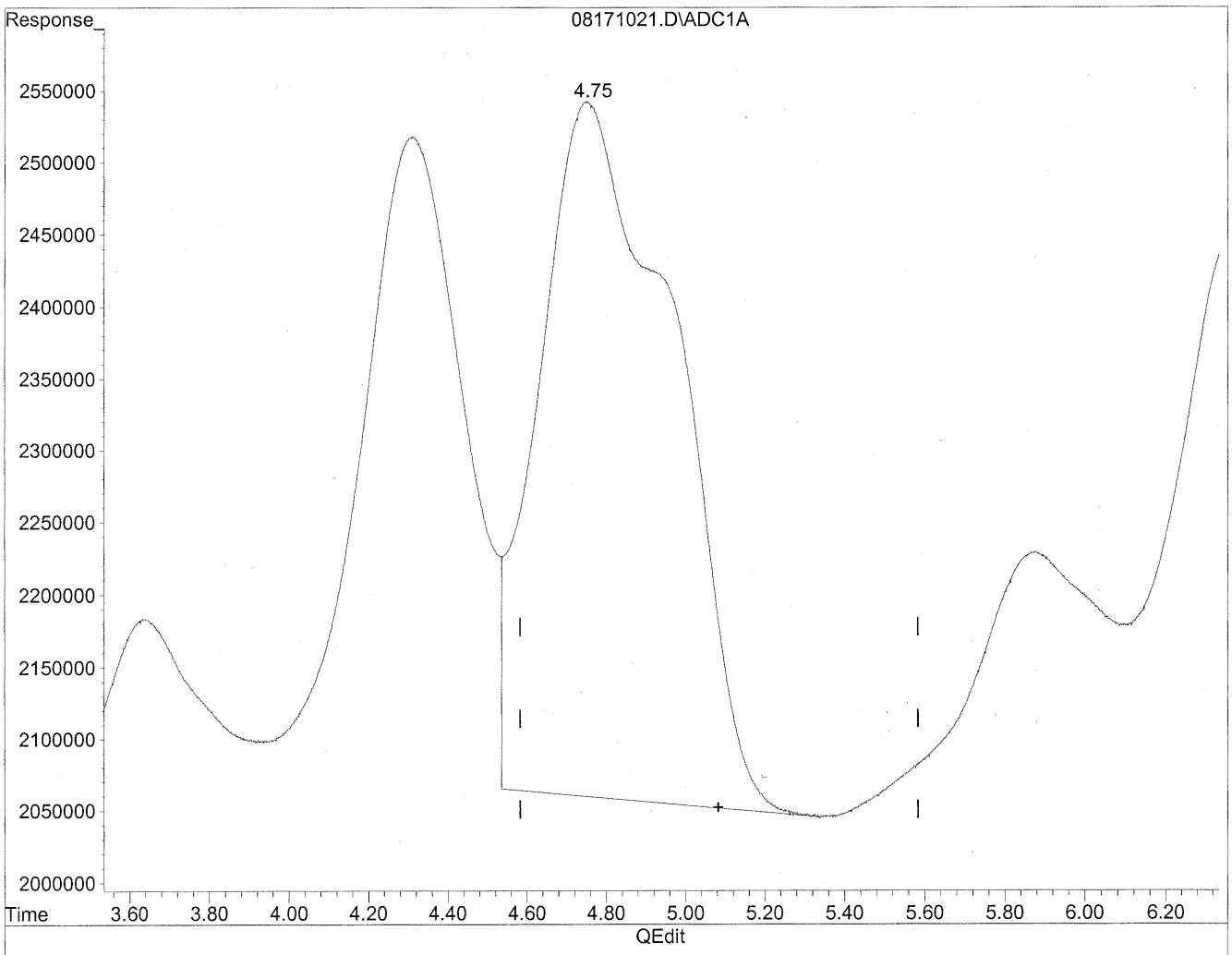
(+) = Expected Retention Time

08171021.D TO110709.M Sat Aug 22 15:24:53 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171021.D Vial: 34
Acq On : 18 Aug 2009 8:54 pm Operator: HC
Sample : P0902772-011 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.75min 1301.661ng/ml

response 114983673

267

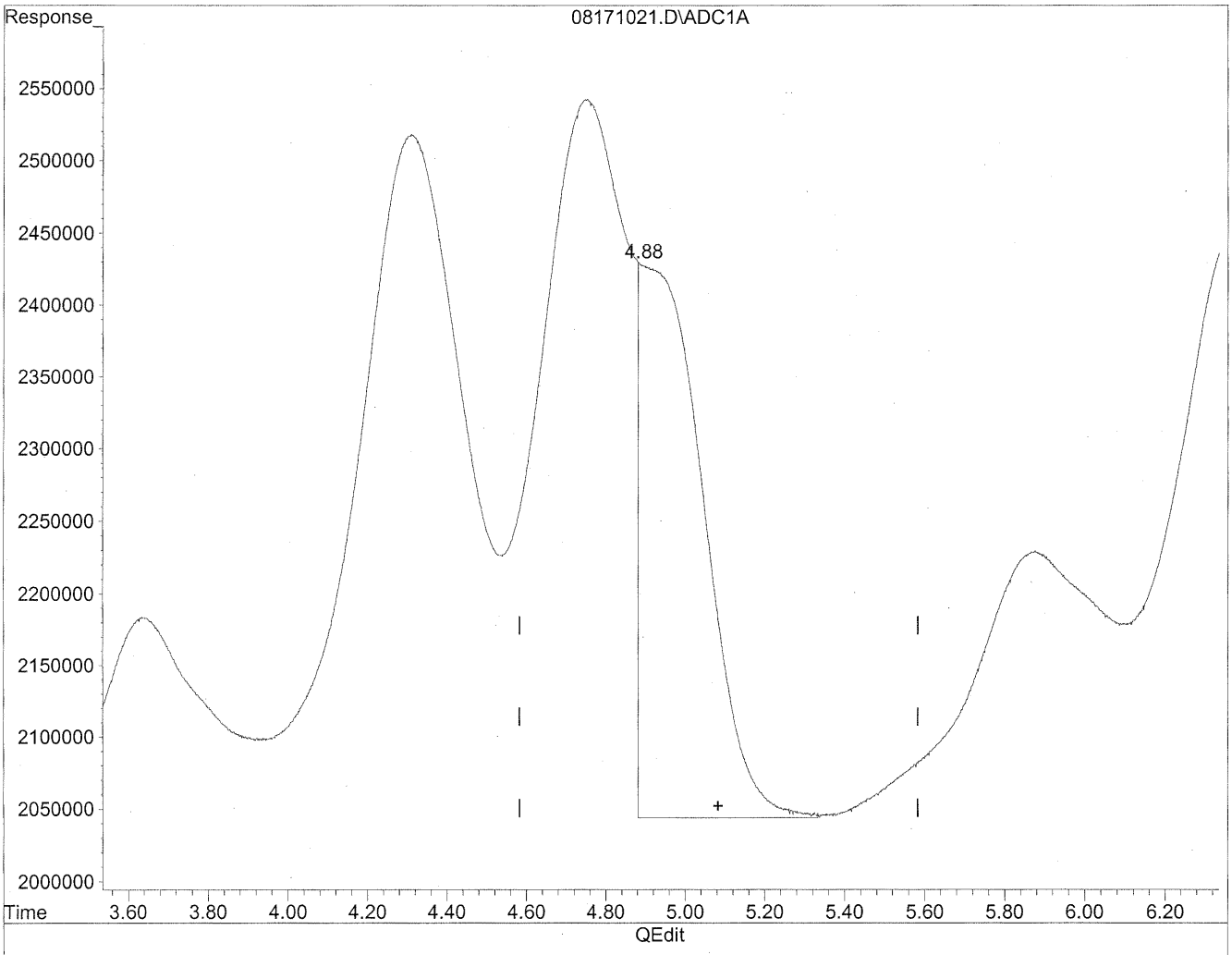
(+) = Expected Retention Time

08171021.D TO110709.M Sat Aug 22 15:25:51 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171021.D Vial: 34
Acq On : 18 Aug 2009 8:54 pm Operator: HC
Sample : P0902772-011 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.88min 478.162ng/ml m

response 42238936

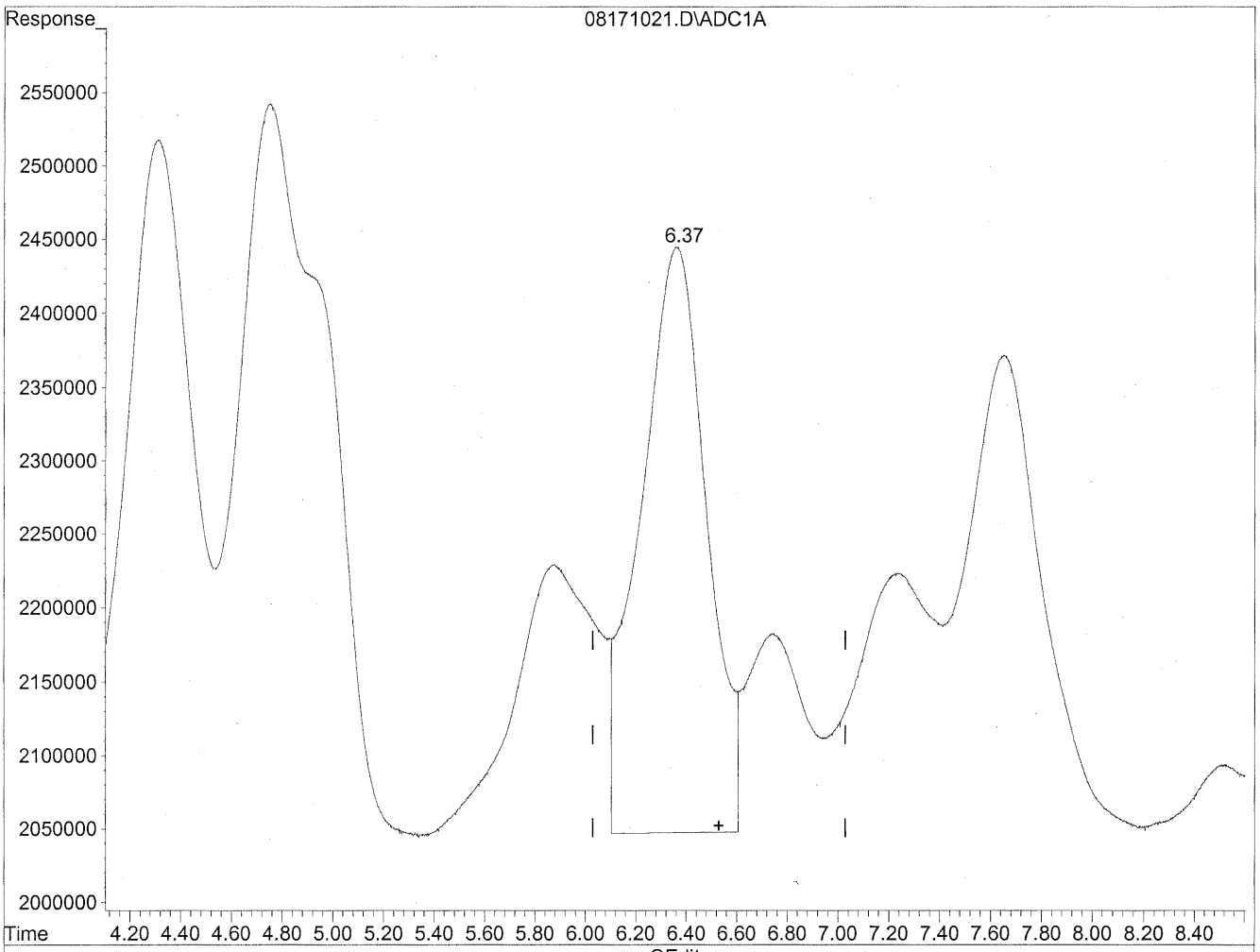
HC 8/22/09
SP
JKP 8/23/09

(+) = Expected Retention Time
08171021.D TO110709.M Sat Aug 22 15:26:10 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171021.D Vial: 34
Acq On : 18 Aug 2009 8:54 pm Operator: HC
Sample : P0902772-011 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.36min 1098.752ng/ml

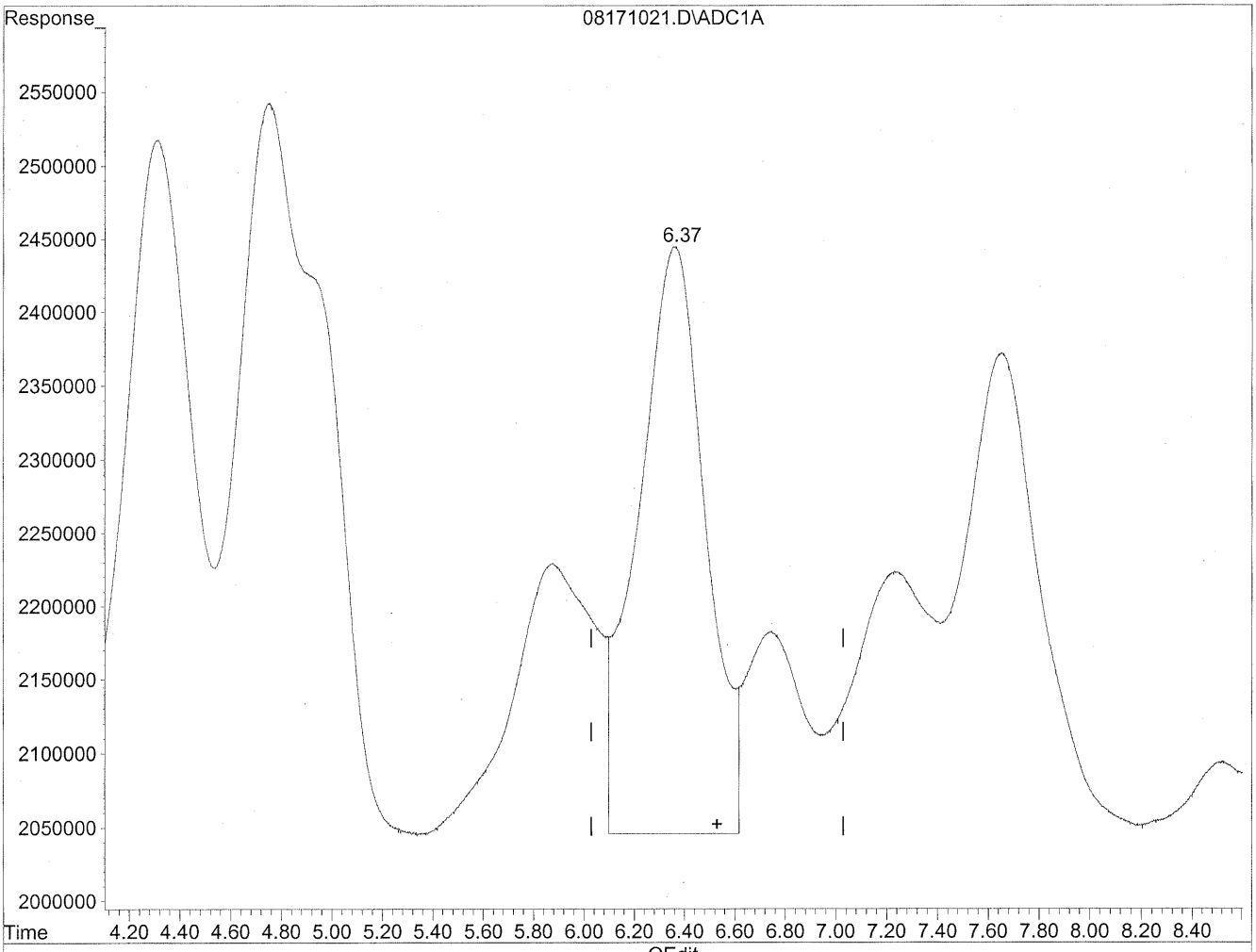
response 72374028

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171021.D Vial: 34
Acq On : 18 Aug 2009 8:54 pm Operator: HC
Sample : P0902772-011 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.37min 1121.396ng/ml m

response 73865582

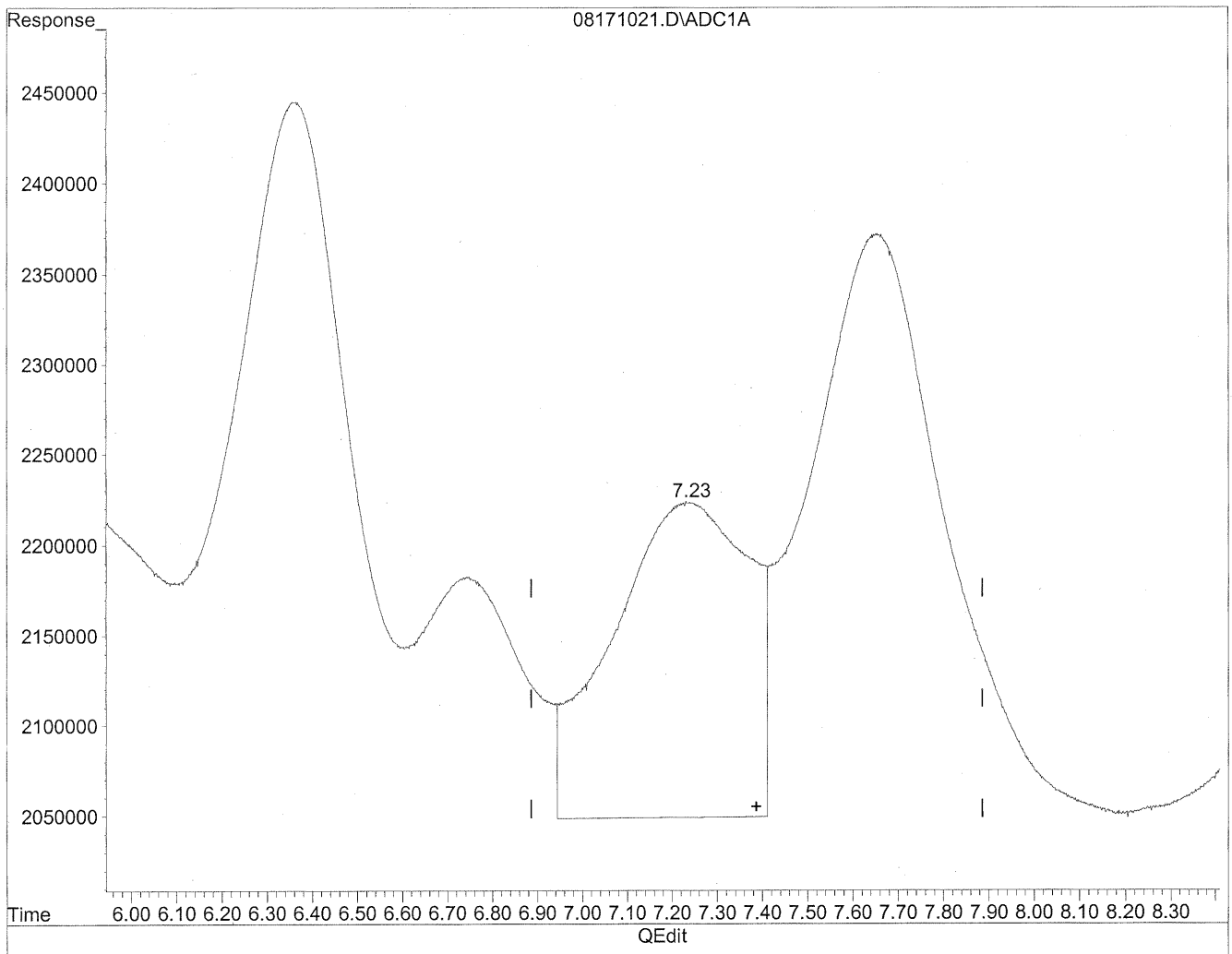
HC 8/22/07
BC
KP 8/23/07

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171021.D Vial: 34
Acq On : 18 Aug 2009 8:54 pm Operator: HC
Sample : P0902772-011 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.24min 469.529ng/ml

response 36741140

271

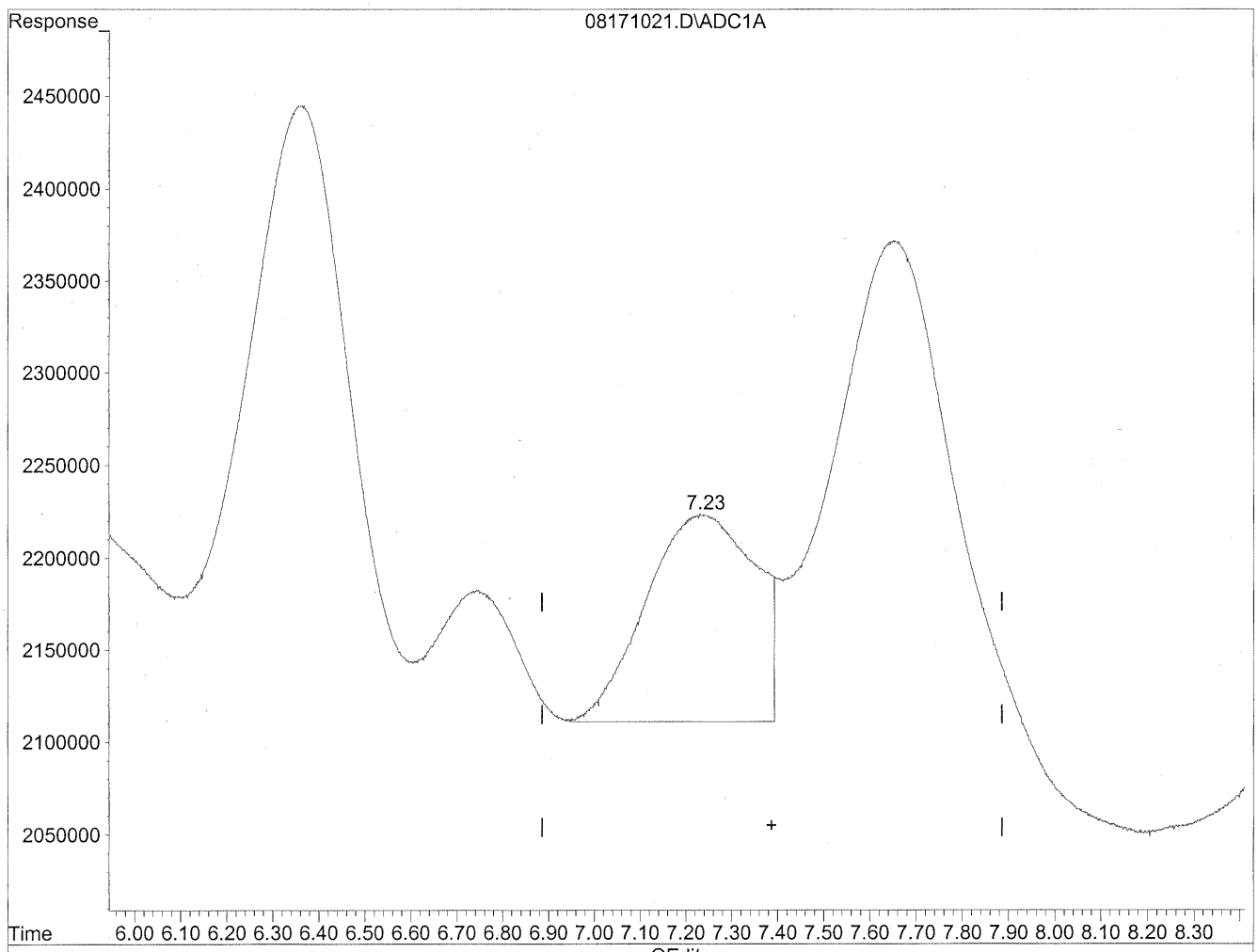
(+) = Expected Retention Time

08171021.D TO110709.M Sat Aug 22 15:26:53 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171021.D Vial: 34
Acq On : 18 Aug 2009 8:54 pm Operator: HC
Sample : P0902772-011 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.23min 237.056ng/ml m

response 18549852

HC 8/22/09
BC

KP 8/23/09

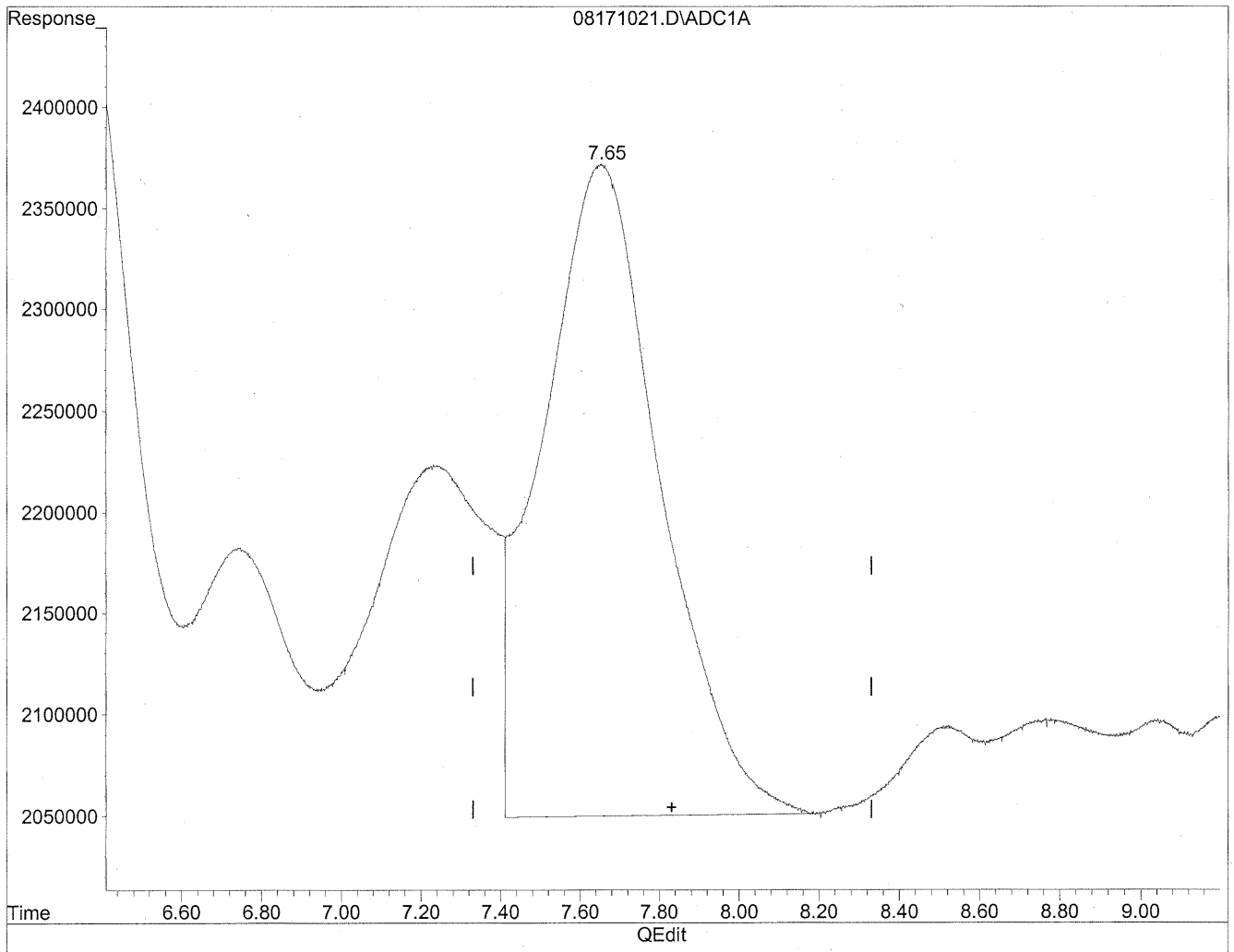
272

(+) = Expected Retention Time
08171021.D TO110709.M Sat Aug 22 15:27:10 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171021.D Vial: 34
Acq On : 18 Aug 2009 8:54 pm Operator: HC
Sample : P0902772-011 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.65min 903.053ng/ml

response 66378923

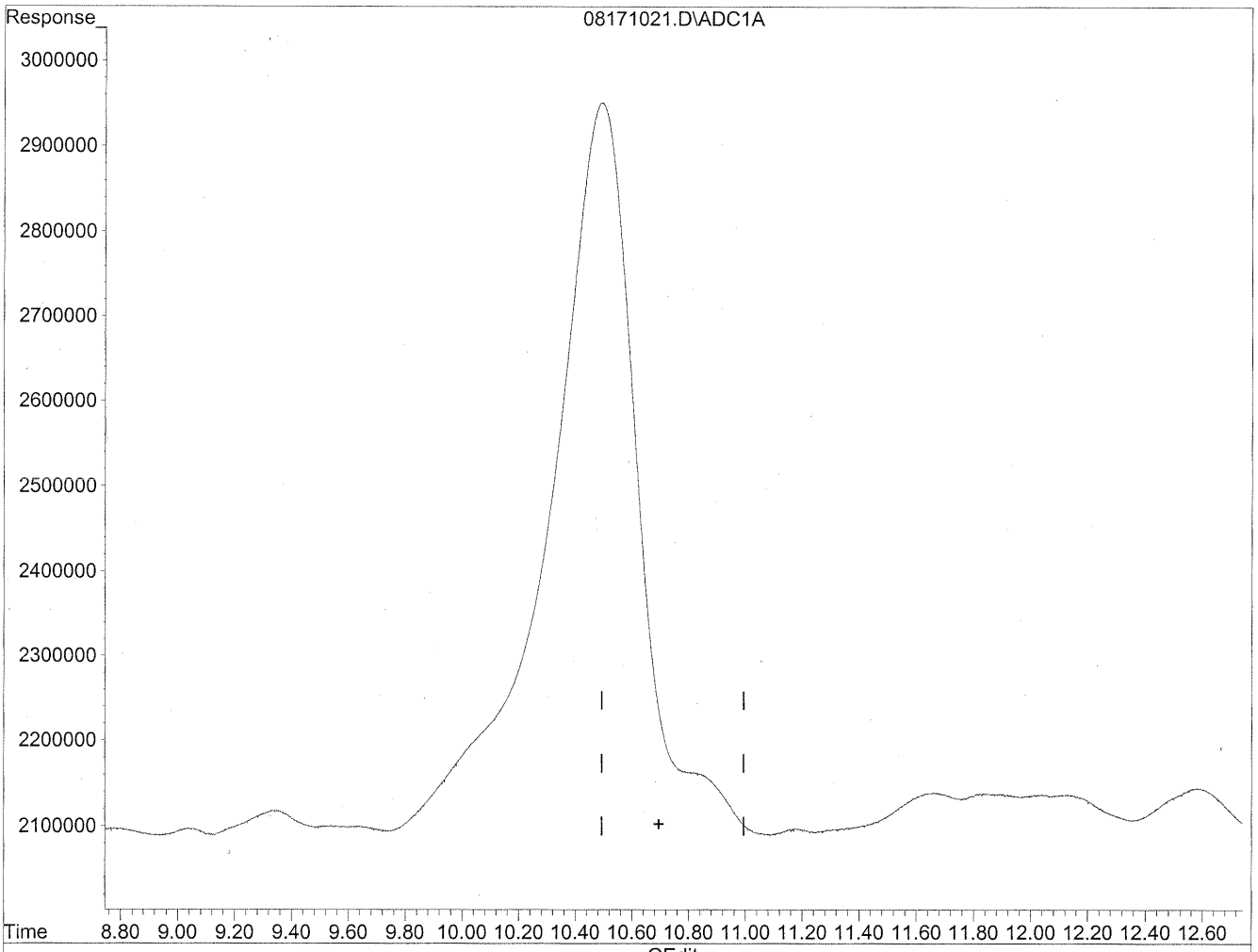
273

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171021.D Vial: 34
Acq On : 18 Aug 2009 8:54 pm Operator: HC
Sample : P0902772-011 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.69min 0.000ng/ml

response 0

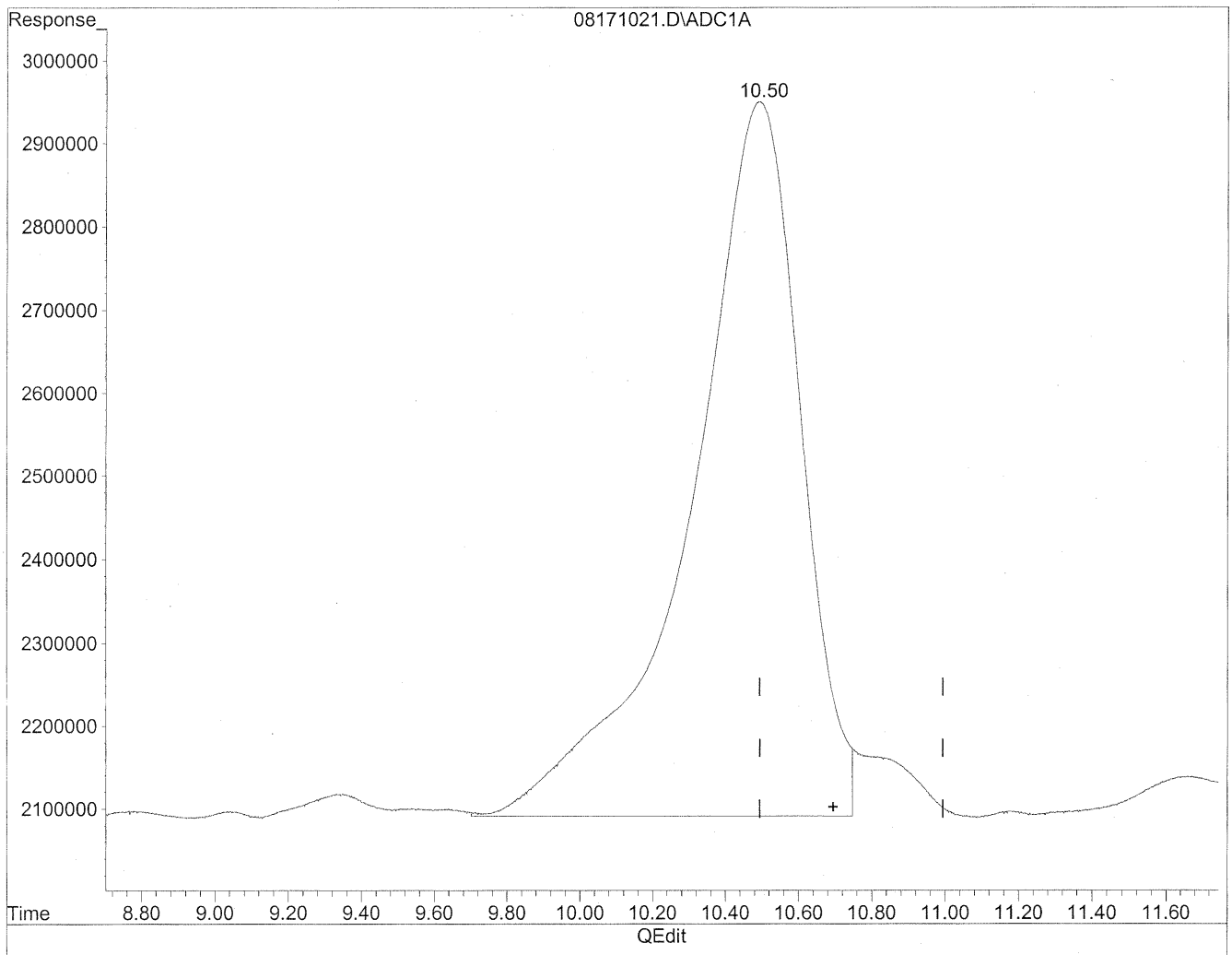
274

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171021.D Vial: 34
Acq On : 18 Aug 2009 8:54 pm Operator: HC
Sample : P0902772-011 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:27 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.50min 2658.393ng/ml m

response 179026219

HC 8/25/09

BNI

MT

KK 8/25/09

275

(+) = Expected Retention Time

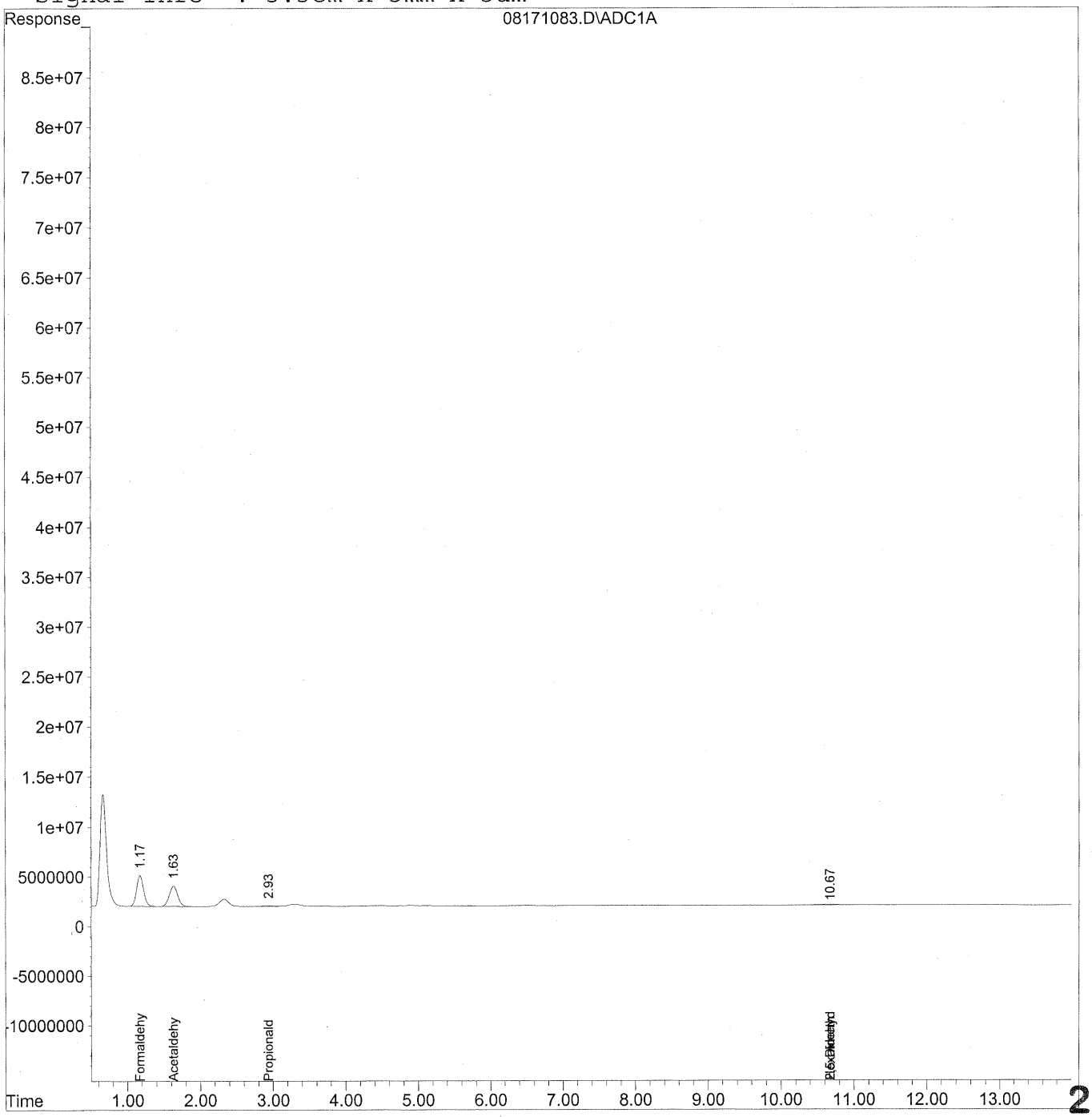
08171021.D TO110709.M Tue Aug 25 07:52:34 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171083.D Vial: 8
Acq On : 19 Aug 2009 12:27 pm Operator: HC
Sample : P0902772-011 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:23 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



276

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171083.D Vial: 8
Acq On : 19 Aug 2009 12:27 pm Operator: HC
Sample : P0902772-011 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:23 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	206160568	1122.993	ng/mlm
2)	Acetaldehyde	1.63	168782410	1203.667	ng/ml
3)	Propionaldehyde	2.93	8028268	75.245	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	10.67	17030812	252.894	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.67	17030812	347.473	ng/ml

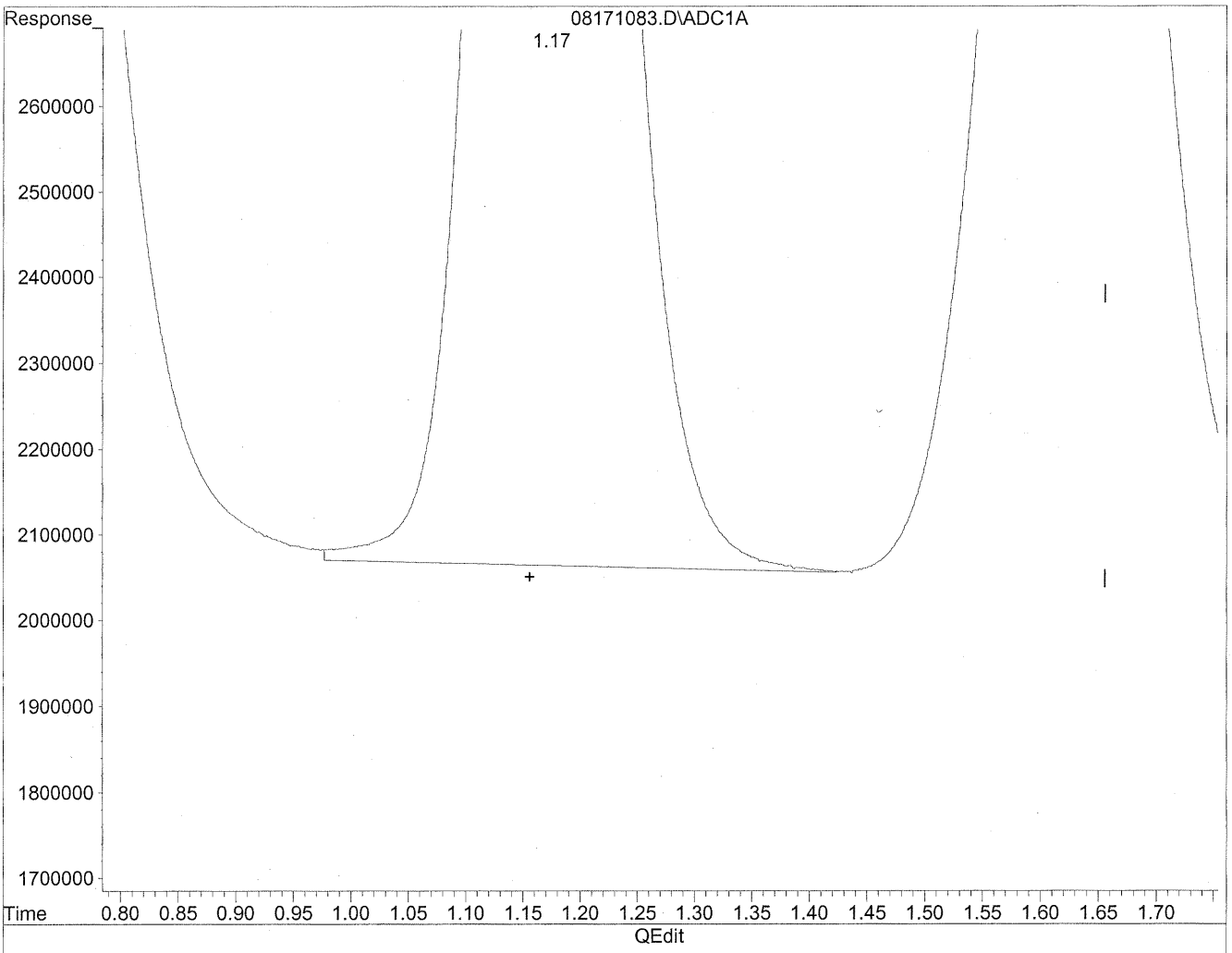
277

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171083.D Vial: 8
Acq On : 19 Aug 2009 12:27 pm Operator: HC
Sample : P0902772-011 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 1132.094ng/ml

response 207831383

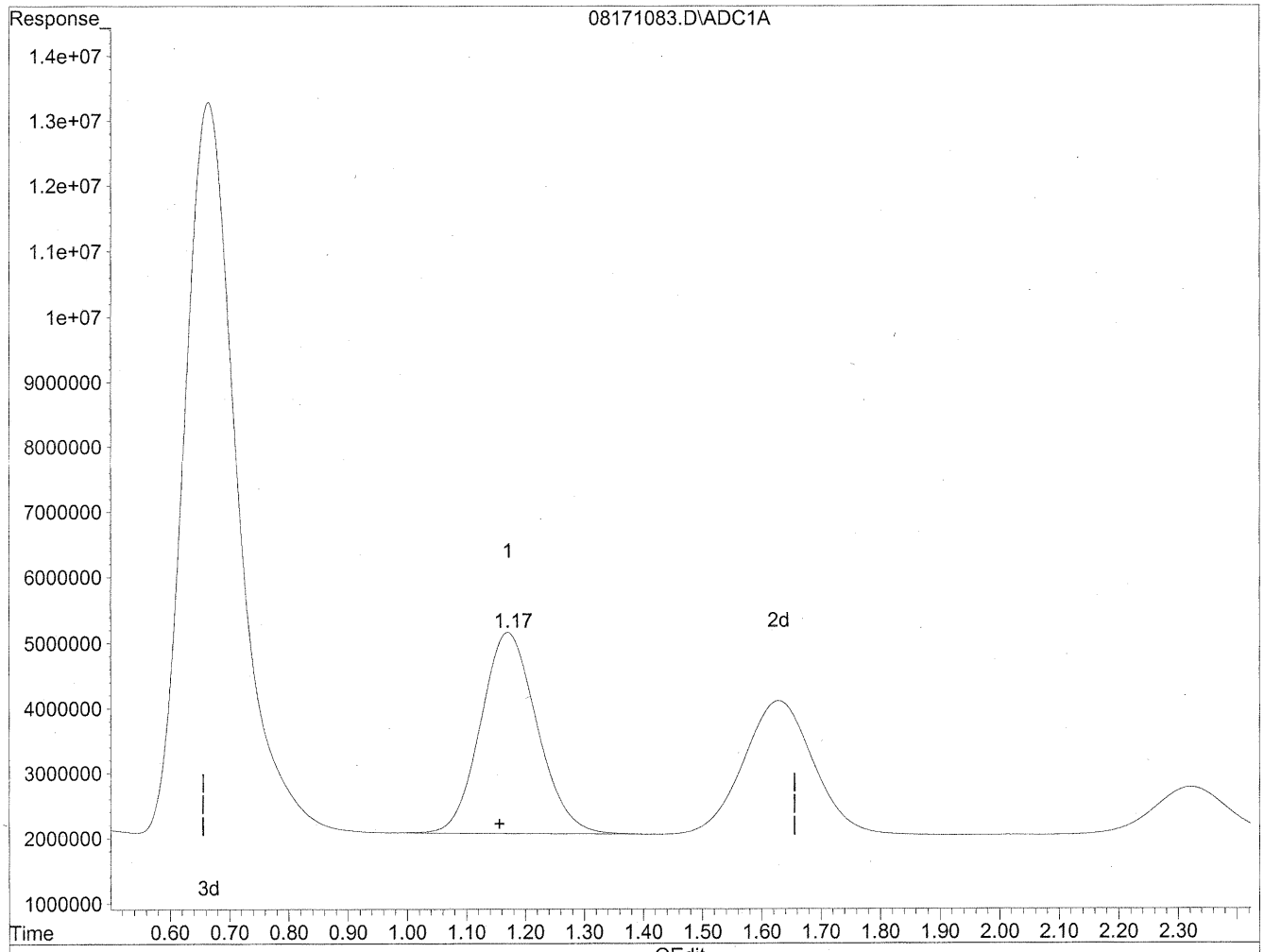
278

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171083.D Vial: 8
Acq On : 19 Aug 2009 12:27 pm Operator: HC
Sample : P0902772-011 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 1122.993ng/ml m

response 206160568

HC 8/22/09
CC
KE 8/23/09

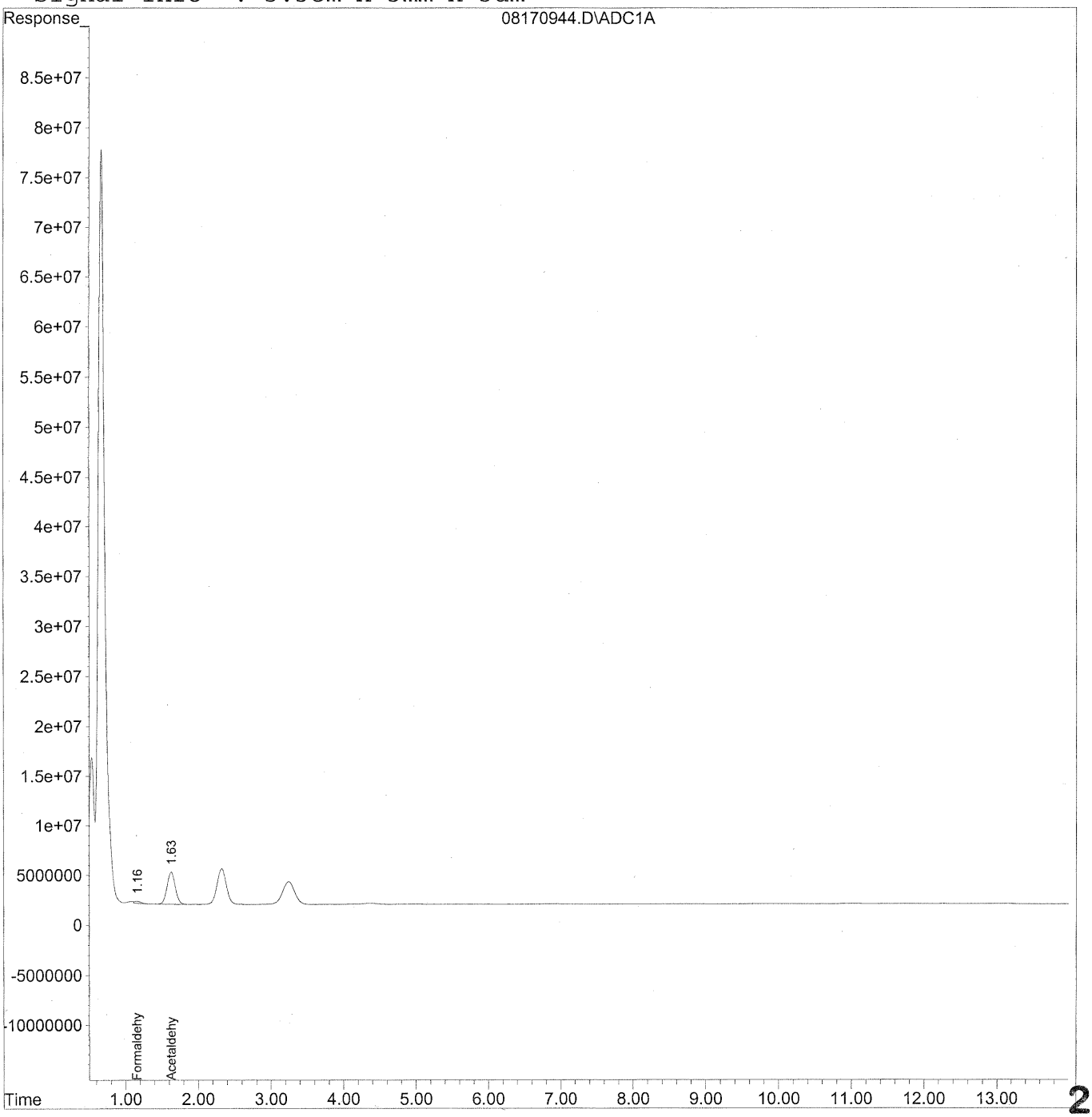
279

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_08\17\08170944.D Vial: 43
Acq On : 18 Aug 2009 1:37 am Operator: HC
Sample : P0902772-011 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:35 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



280

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170944.D Vial: 43
Acq On : 18 Aug 2009 1:37 am Operator: HC
Sample : P0902772-011 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:35 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

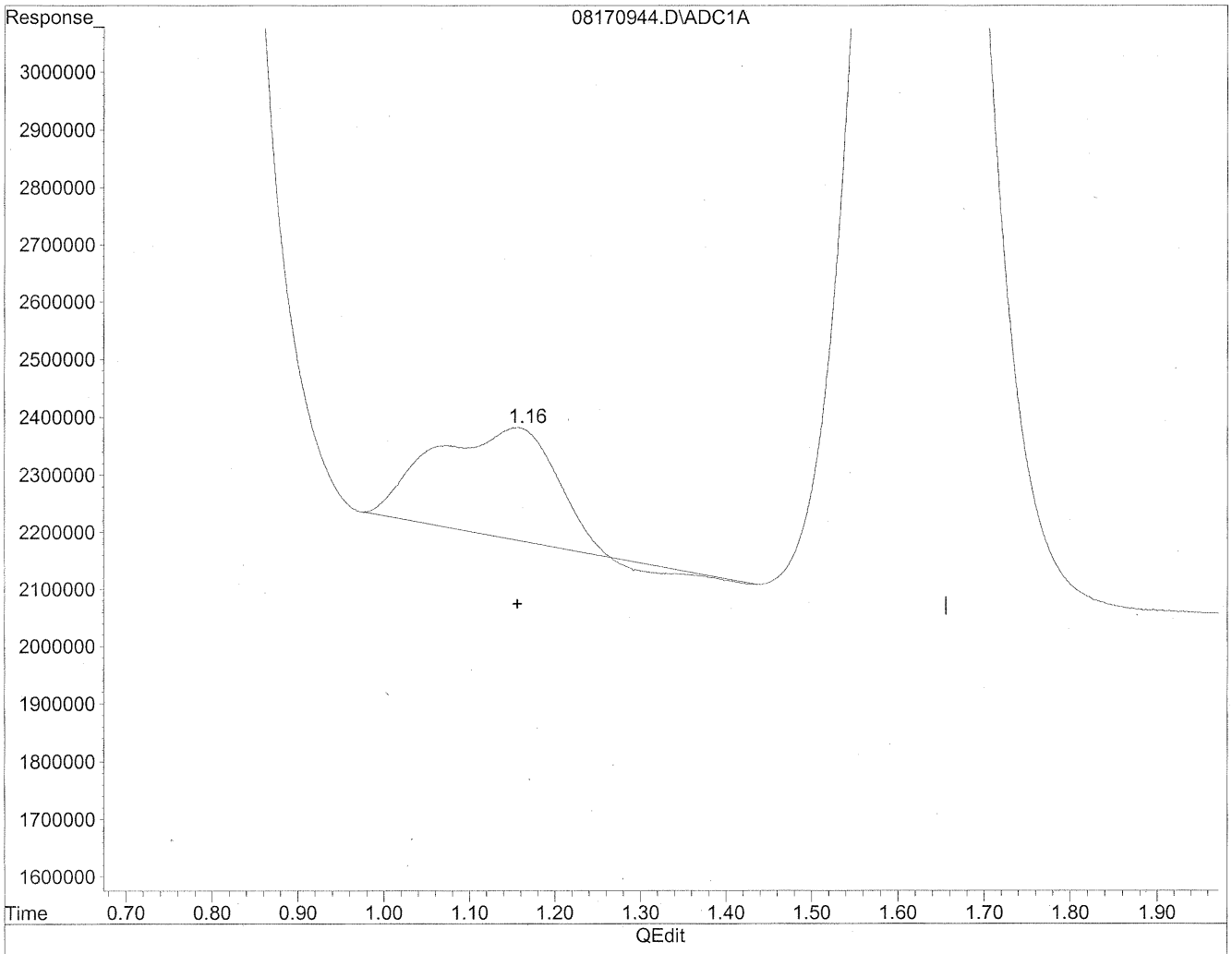
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	12236295	66.653	ng/mlm
2)	Acetaldehyde	1.63	254498309	1814.948	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170944.D Vial: 43
Acq On : 18 Aug 2009 1:37 am Operator: HC
Sample : P0902772-011 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:34 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.16min 100.249ng/ml

response 18403831

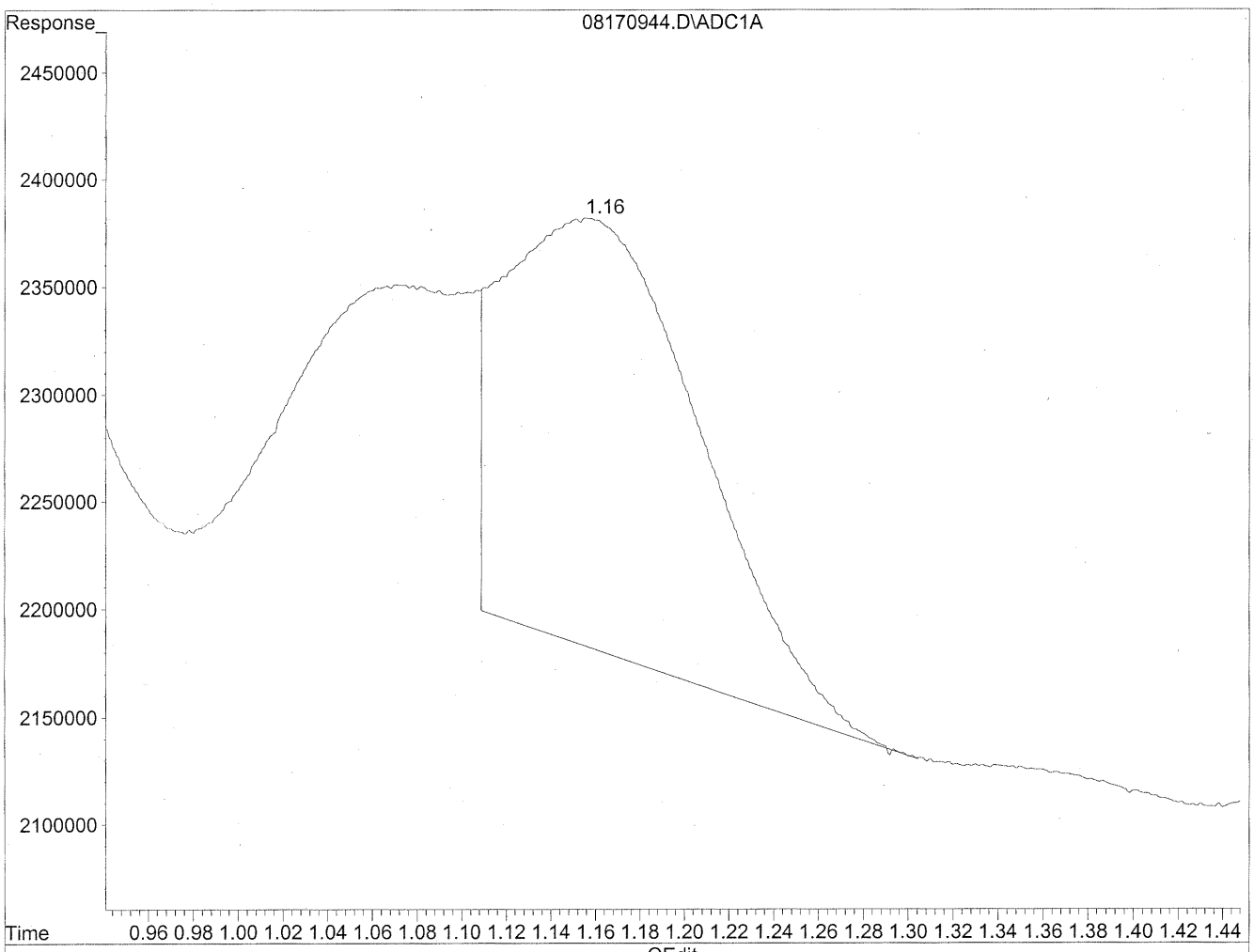
(+) = Expected Retention Time

08170944.D TO110709.M Sat Aug 22 12:35:09 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170944.D Vial: 43
Acq On : 18 Aug 2009 1:37 am Operator: HC
Sample : P0902772-011 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:34 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.16min 66.653ng/ml m

response 12236295

HC 8/21/09 SP

KP 8/23/09

283

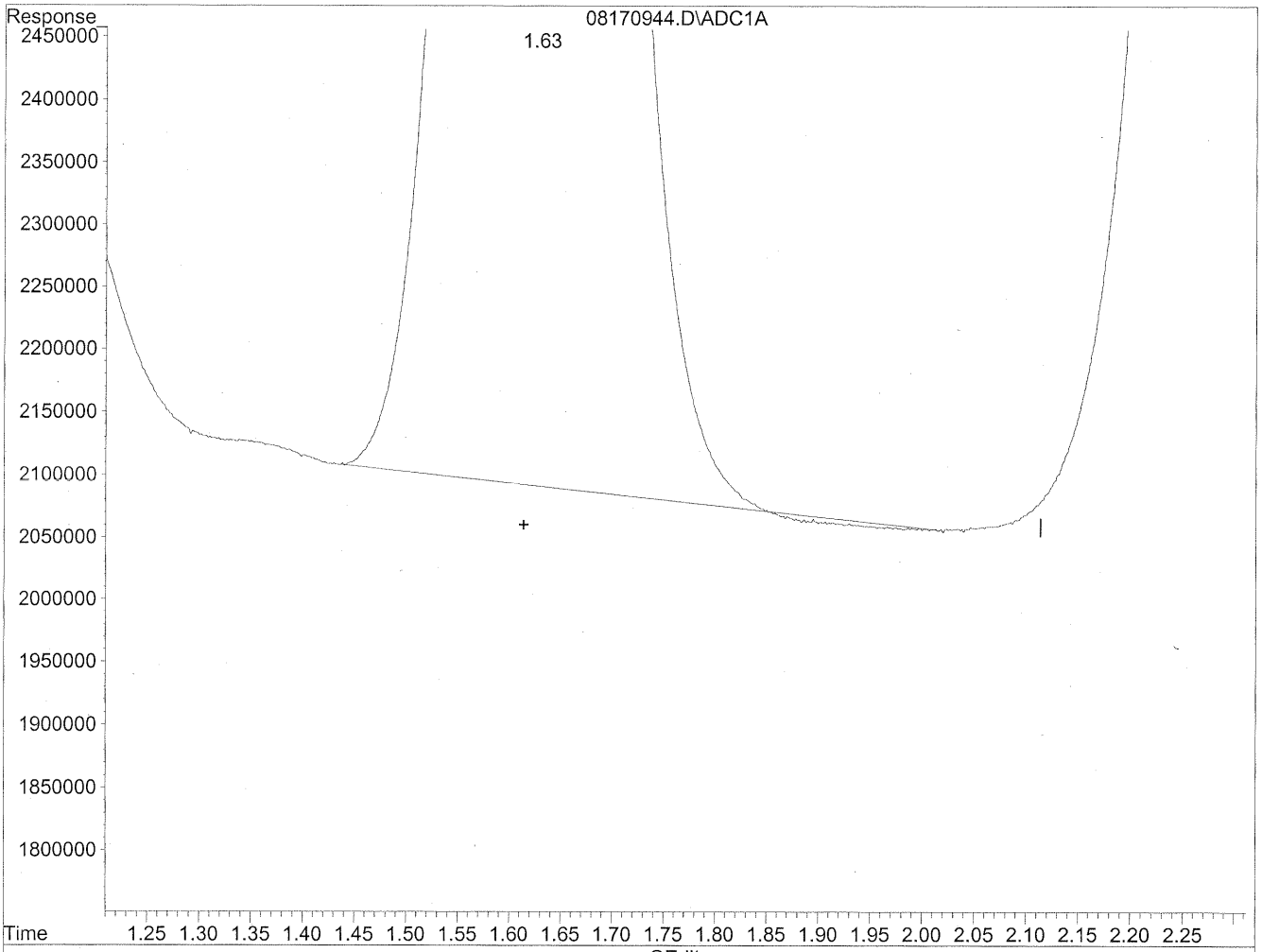
(+) = Expected Retention Time

08170944.D TO110709.M Sat Aug 22 12:35:17 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170944.D Vial: 43
Acq On : 18 Aug 2009 1:37 am Operator: HC
Sample : P0902772-011 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:34 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.63min 1808.861ng/ml

response 253644783

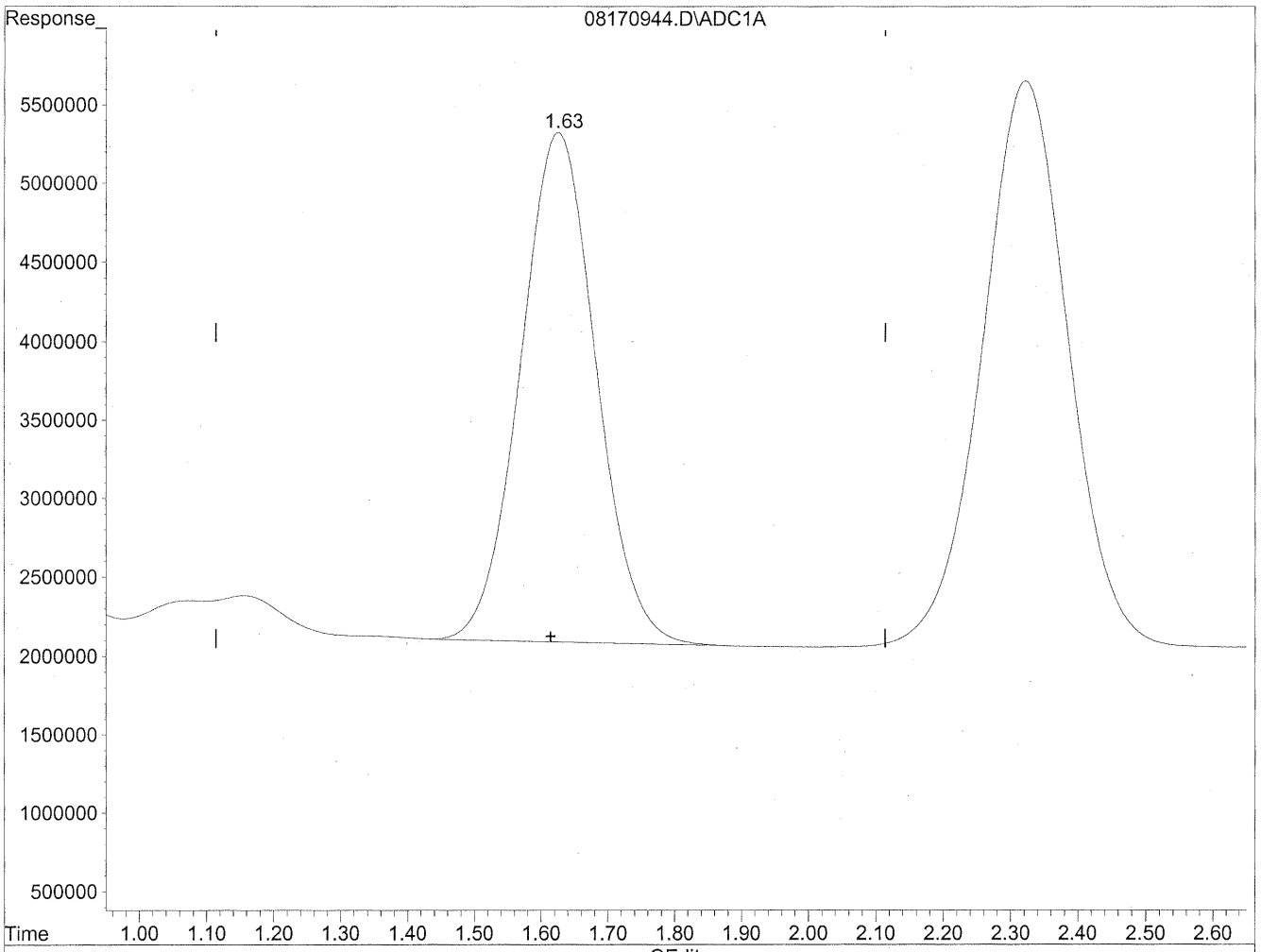
(+) = Expected Retention Time

08170944.D TO110709.M Sat Aug 22 12:35:26 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170944.D Vial: 43
Acq On : 18 Aug 2009 1:37 am Operator: HC
Sample : P0902772-011 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:34 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.63min 1814.948ng/ml m

response 254498309

HC 8/21/09 LC

KP 8/23/09

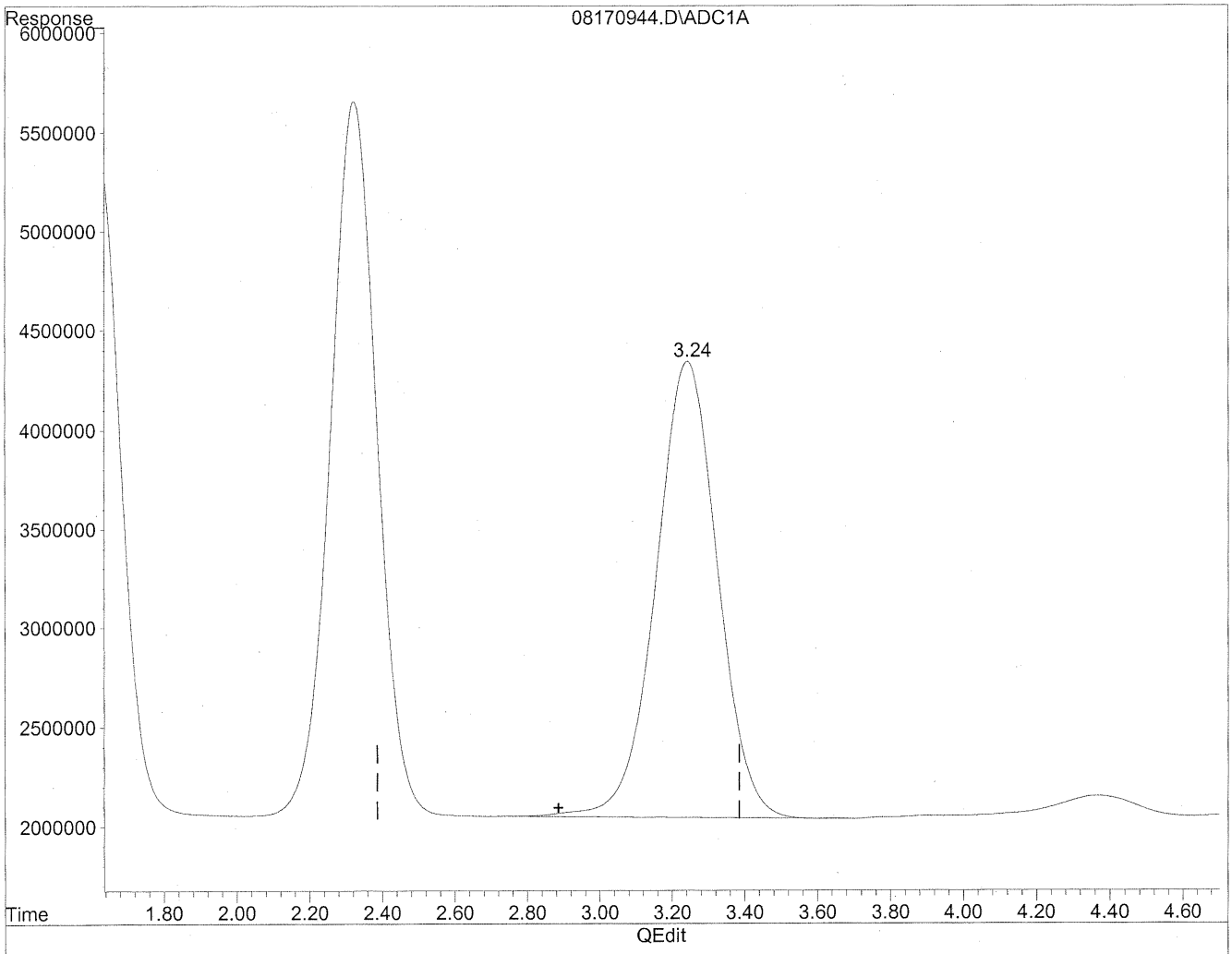
285

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170944.D Vial: 43
Acq On : 18 Aug 2009 1:37 am Operator: HC
Sample : P0902772-011 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:34 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.24min 2549.663ng/ml

response 272036912

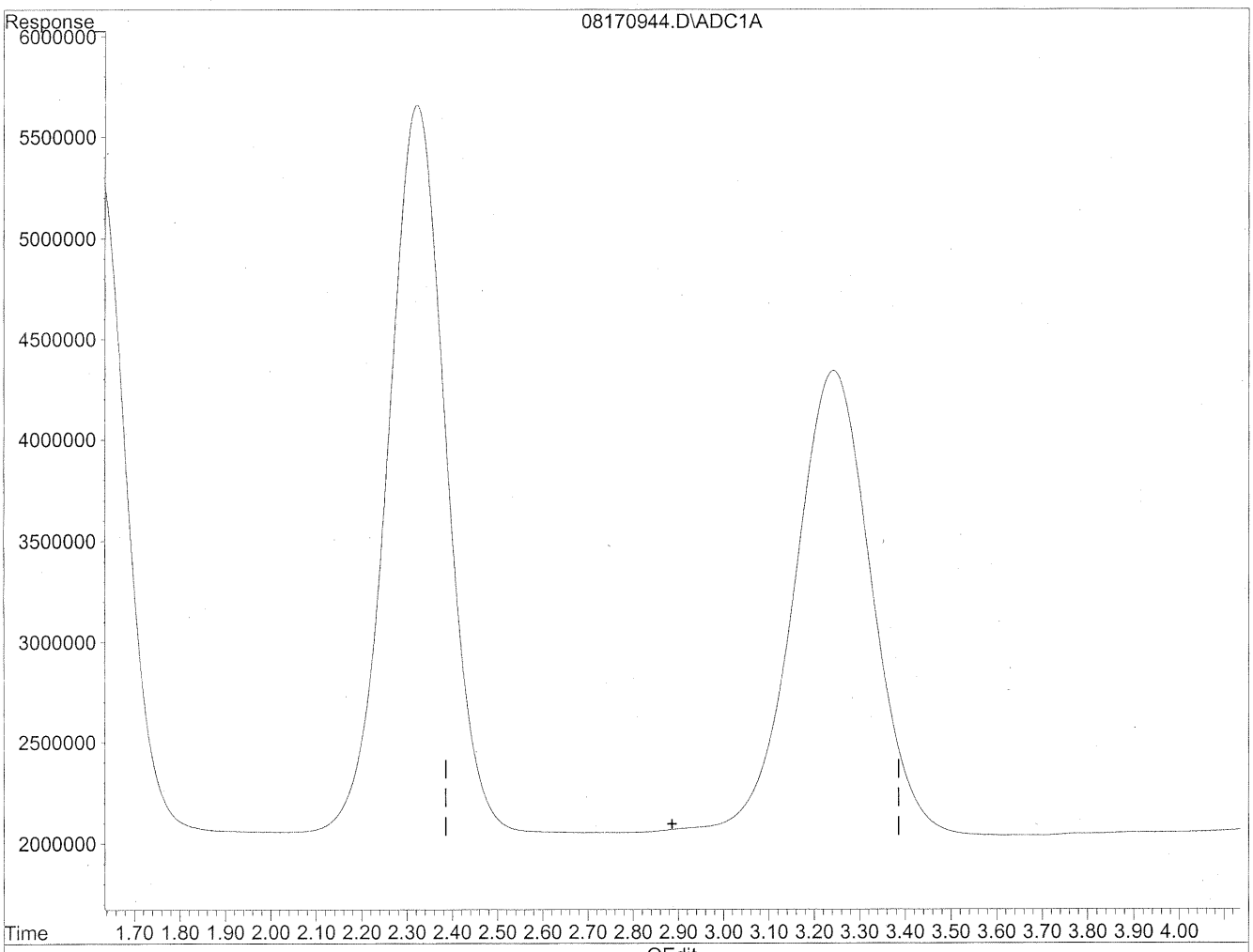
(+) = Expected Retention Time

08170944.D TO110709.M Sat Aug 22 12:35:42 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170944.D Vial: 43
Acq On : 18 Aug 2009 1:37 am Operator: HC
Sample : P0902772-011 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:35 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

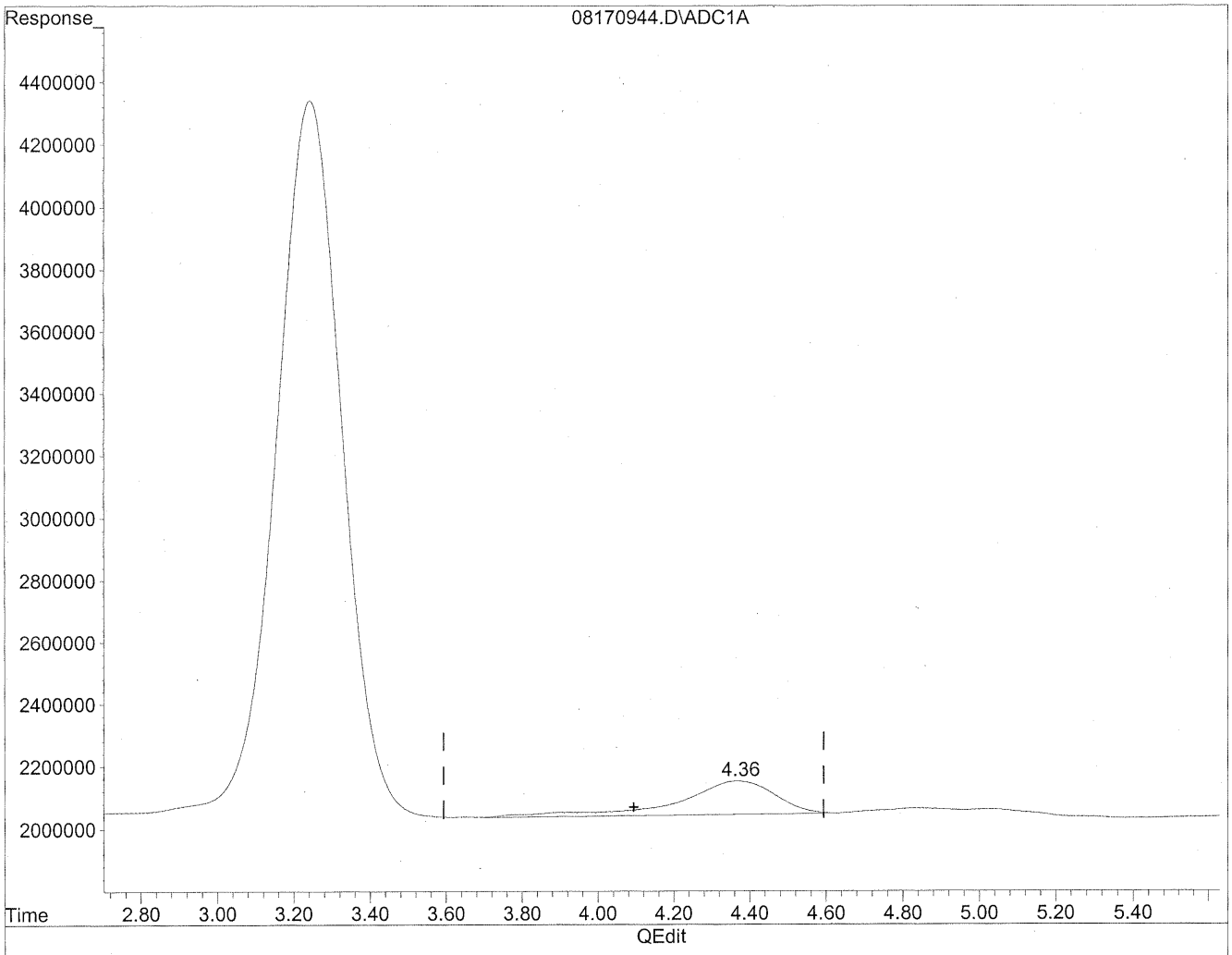
response 0

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170944.D Vial: 43
Acq On : 18 Aug 2009 1:37 am Operator: HC
Sample : P0902772-011 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:34 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.37min 192.409ng/ml

response 18743520

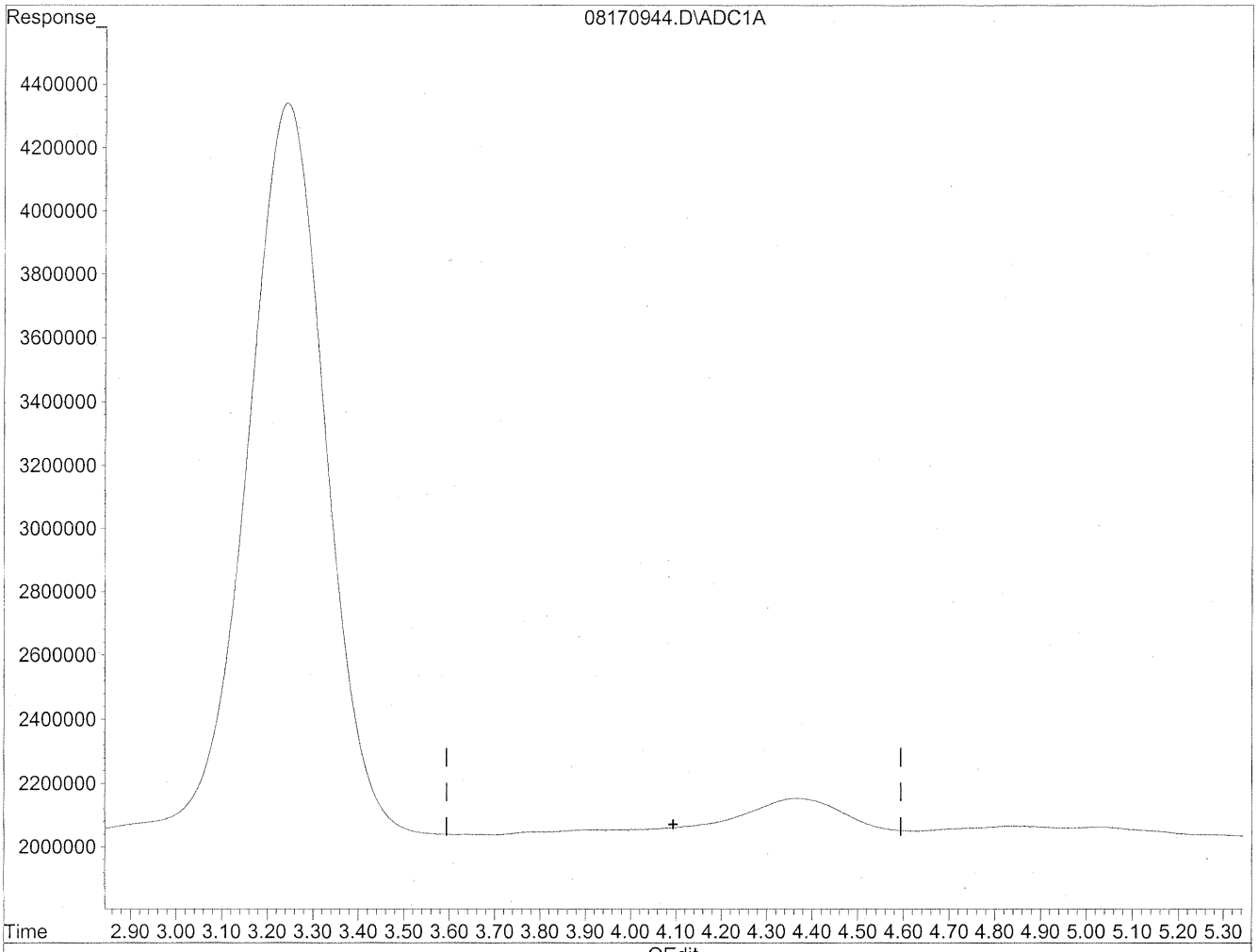
(+) = Expected Retention Time

08170944.D TO110709.M Sat Aug 22 12:35:53 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170944.D Vial: 43
Acq On : 18 Aug 2009 1:37 am Operator: HC
Sample : P0902772-011 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:35 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/25/09 MP

HC 8/25/09

289

(+) = Expected Retention Time

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **100436**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P0902772-012

Test Code: EPA Method TO-11A
Instrument ID: Waters LC Module I Plus/UV_Vis 360/LC1
Analyst: Hani Cherazaie
Sampling Media: Silica Gel DNPH Tube
Test Notes: **BC**

Date Collected: 8/10/09
Date Received: 8/12/09
Date Analyzed: 8/18 - 8/19/09
Desorption Volume: 1.0 ml
Volume Sampled: 103.7 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**12,000**	**110**	0.96	**91**	0.79	
75-07-0	**Acetaldehyde**	**14,000**	**130**	0.96	**74**	0.54	**BT**
123-38-6	**Propionaldehyde**	**860**	**8.3**	0.96	**3.5**	0.41	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.96	ND	0.34	
123-72-8	**Butyraldehyde**	**450**	**4.3**	0.96	**1.5**	0.33	
100-52-7	**Benzaldehyde**	**1,300**	**13**	0.96	**3.0**	0.22	
590-86-3	**Isovaleraldehyde**	**340**	**3.3**	0.96	**0.93**	0.27	
110-62-3	**Valeraldehyde**	**1,200**	**11**	0.96	**3.2**	0.27	
529-20-4	o-Tolualdehyde	< 100	ND	0.96	ND	0.20	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	1.9	ND	0.39	
66-25-1	**n-Hexaldehyde**	**3,000**	**29**	0.96	**7.0**	0.24	
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	ND	0.96	ND	0.18	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
BT = Results indicated possible breakthrough; back section > 10% front section.

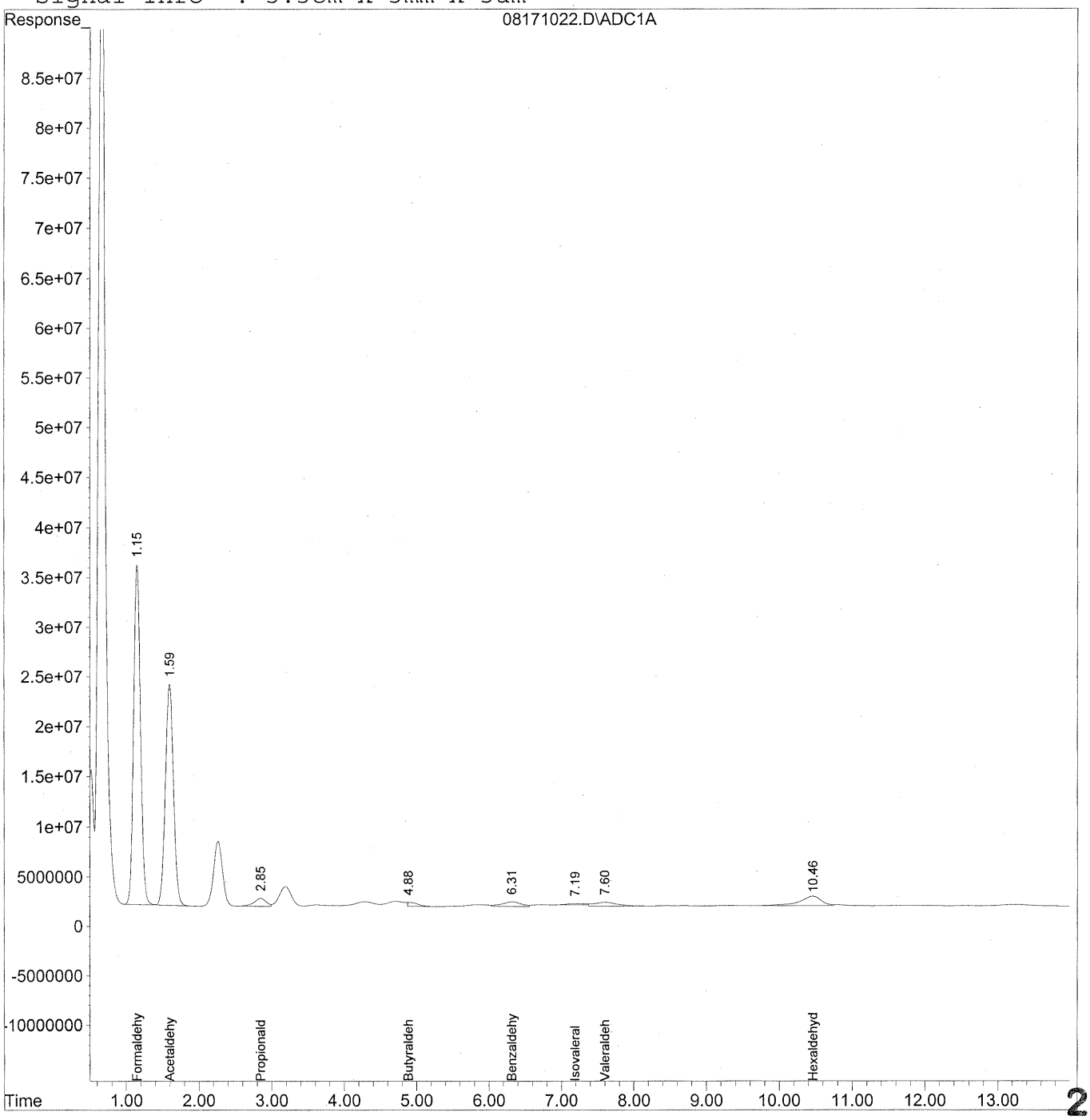
Verified By: _____ Date: 8/25/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:32 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



291

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:32 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

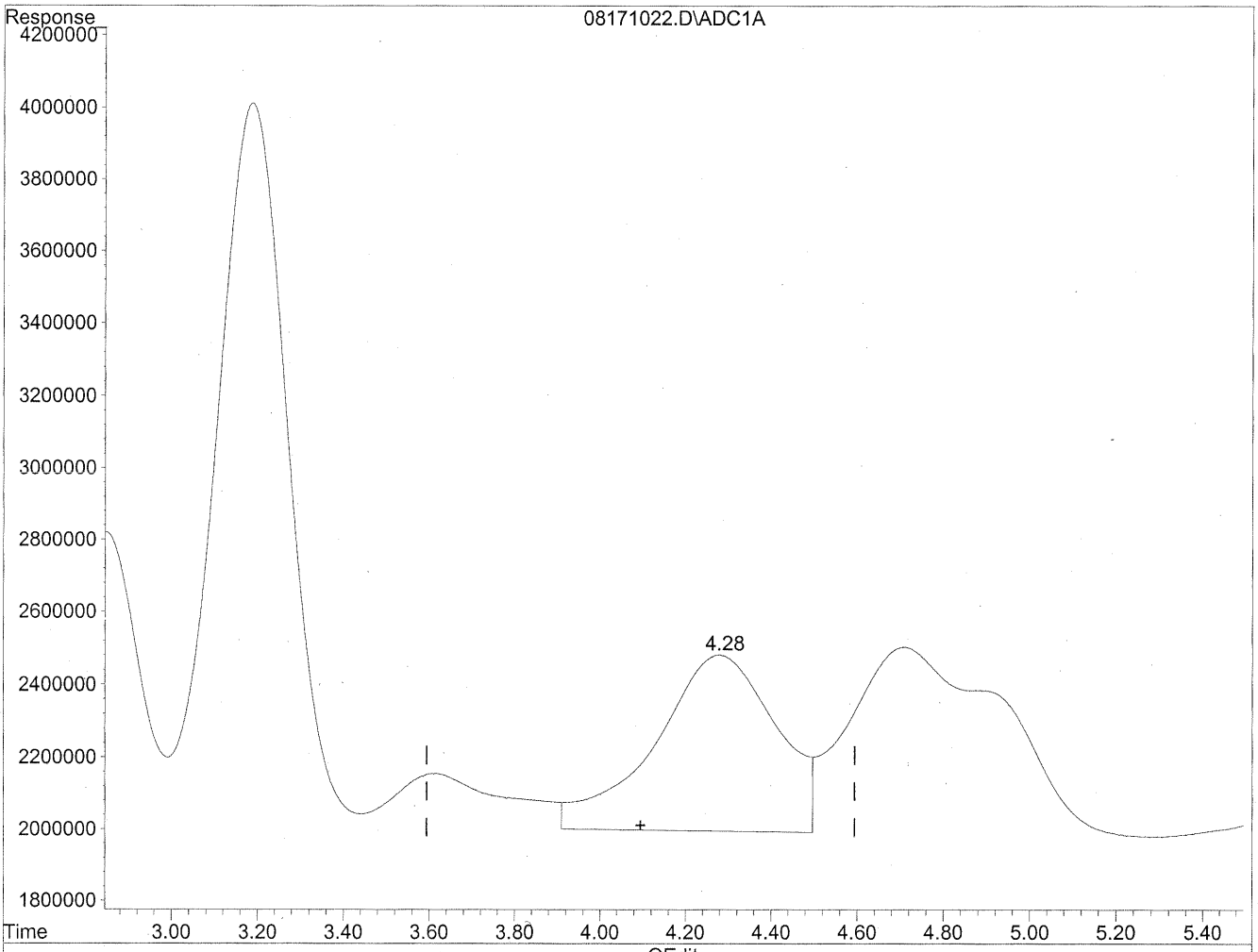
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.15	2205673173	12014.688	ng/ml
2)	Acetaldehyde	1.59	1727519348	12319.756	ng/ml
3)	Propionaldehyde	2.85	92097675	863.185	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	4.88	39552170	447.746	ng/mlm
6)	Benzaldehyde	6.31	88088777	1337.327	ng/mlm
7)	Isovaleraldehyde	7.19	26563770	339.469	ng/mlm
8)	Valeraldehyde	7.60	85516374	1163.409	ng/mlm
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	10.46	201344667	2989.804	ng/mlm
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/mld

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:28 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.28min 985.802ng/ml

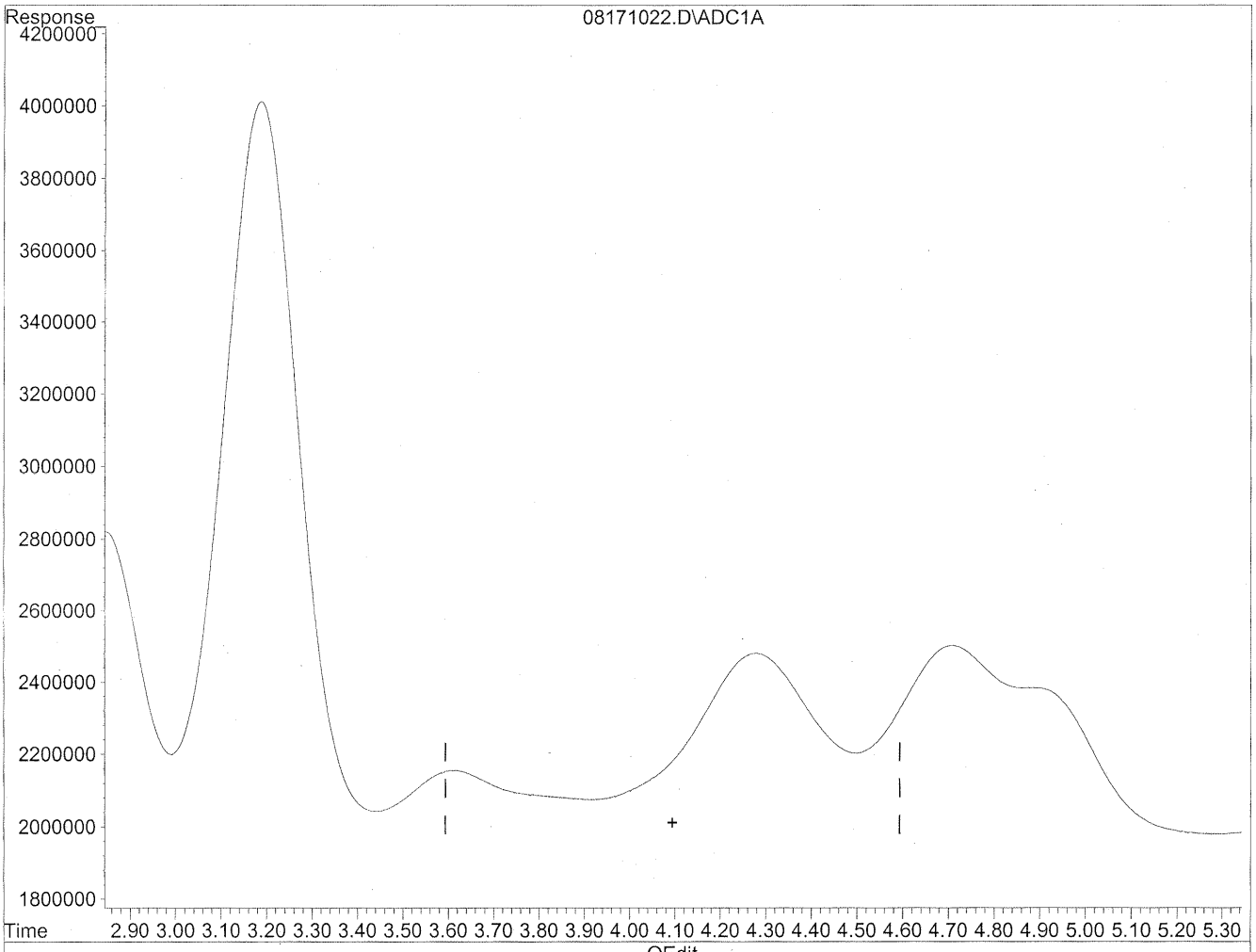
response 96032069

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:32 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

dc 8/28/09 mp

8/25/09

294

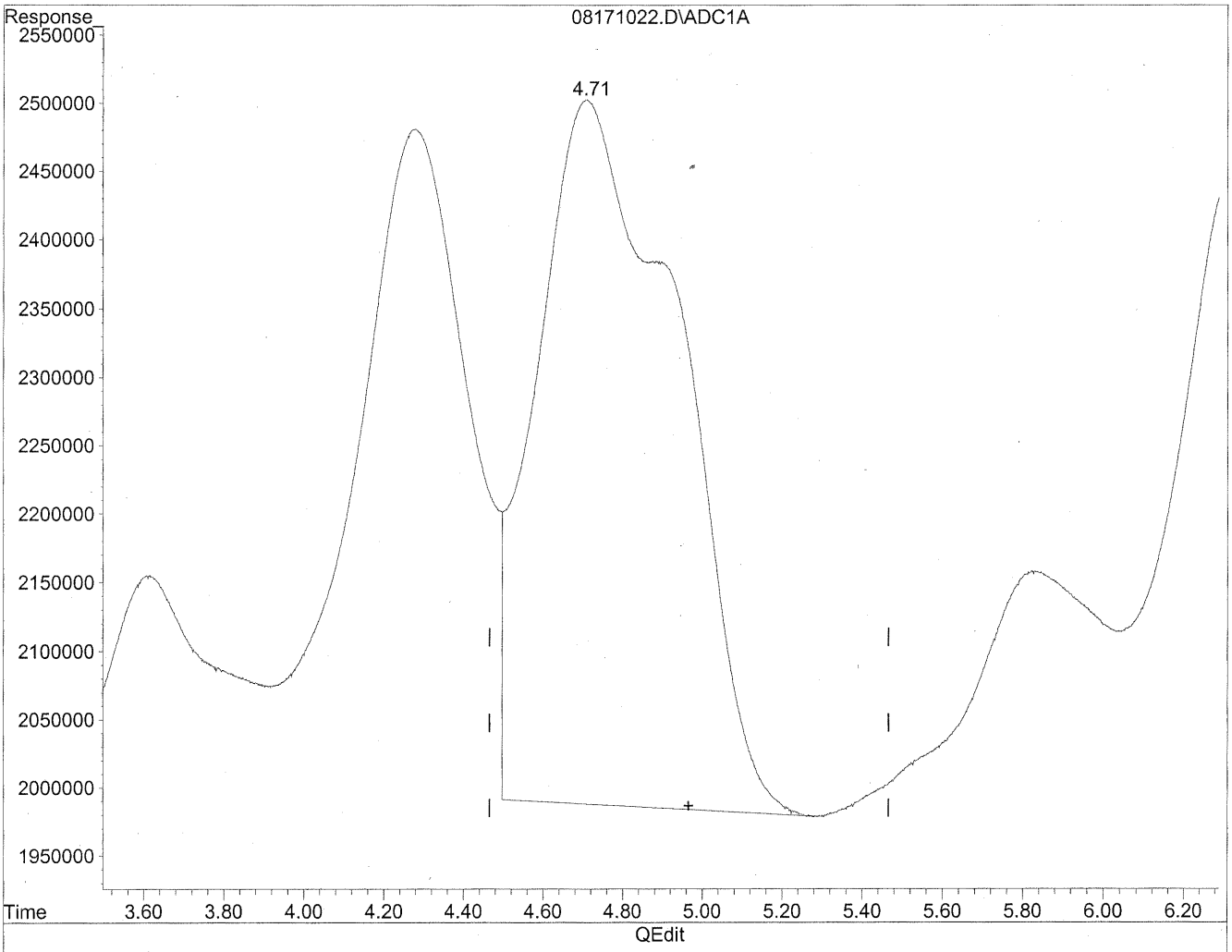
(+) = Expected Retention Time

08171022.D TO110709.M Tue Aug 25 07:53:43 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:28 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.71min 1449.927ng/ml

response 128080934

295

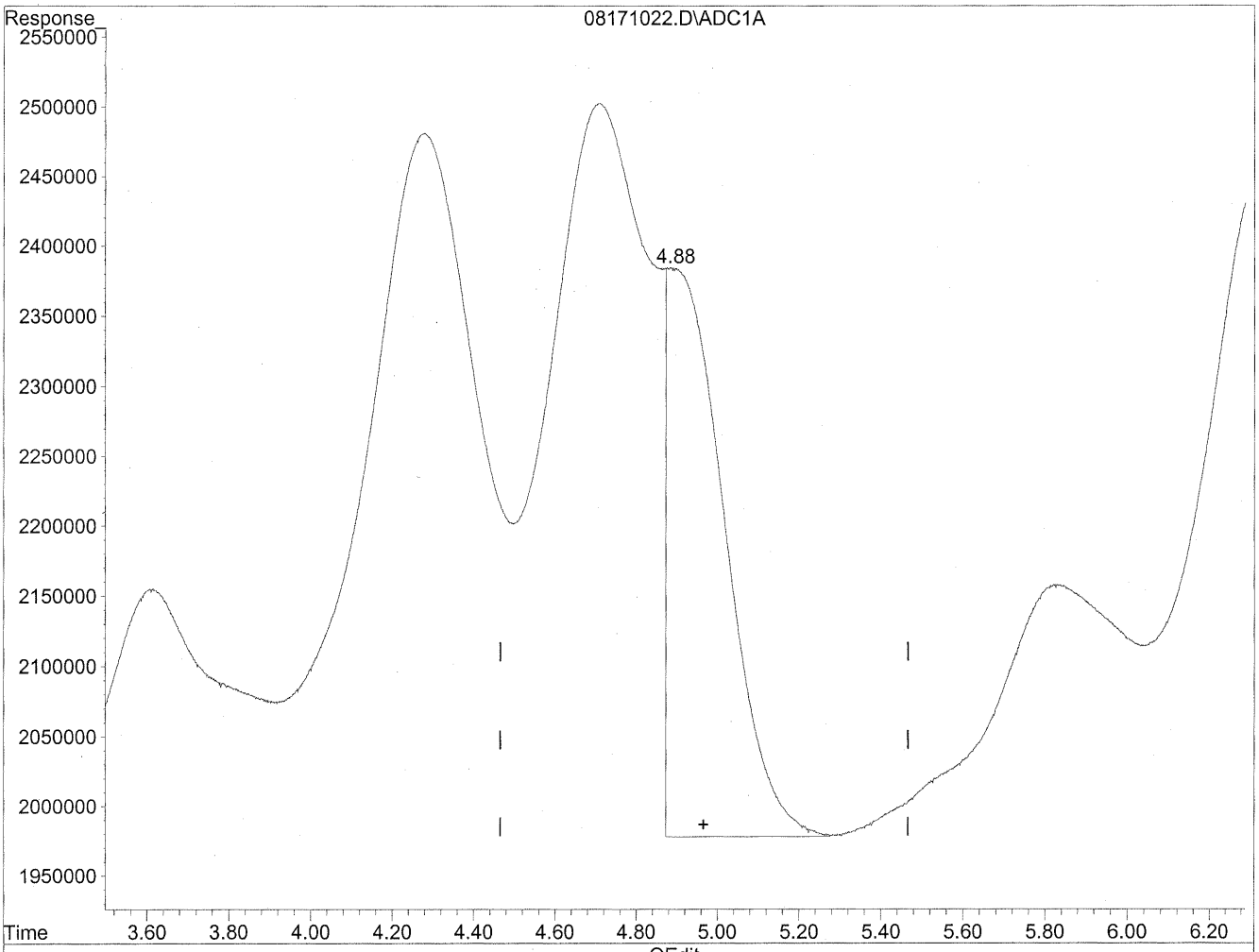
(+) = Expected Retention Time

08171022.D TO110709.M Sat Aug 22 15:29:46 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:28 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.88min 447.746ng/ml m

response 39552170

HC 8/22/09
SP

KR 8/23/09

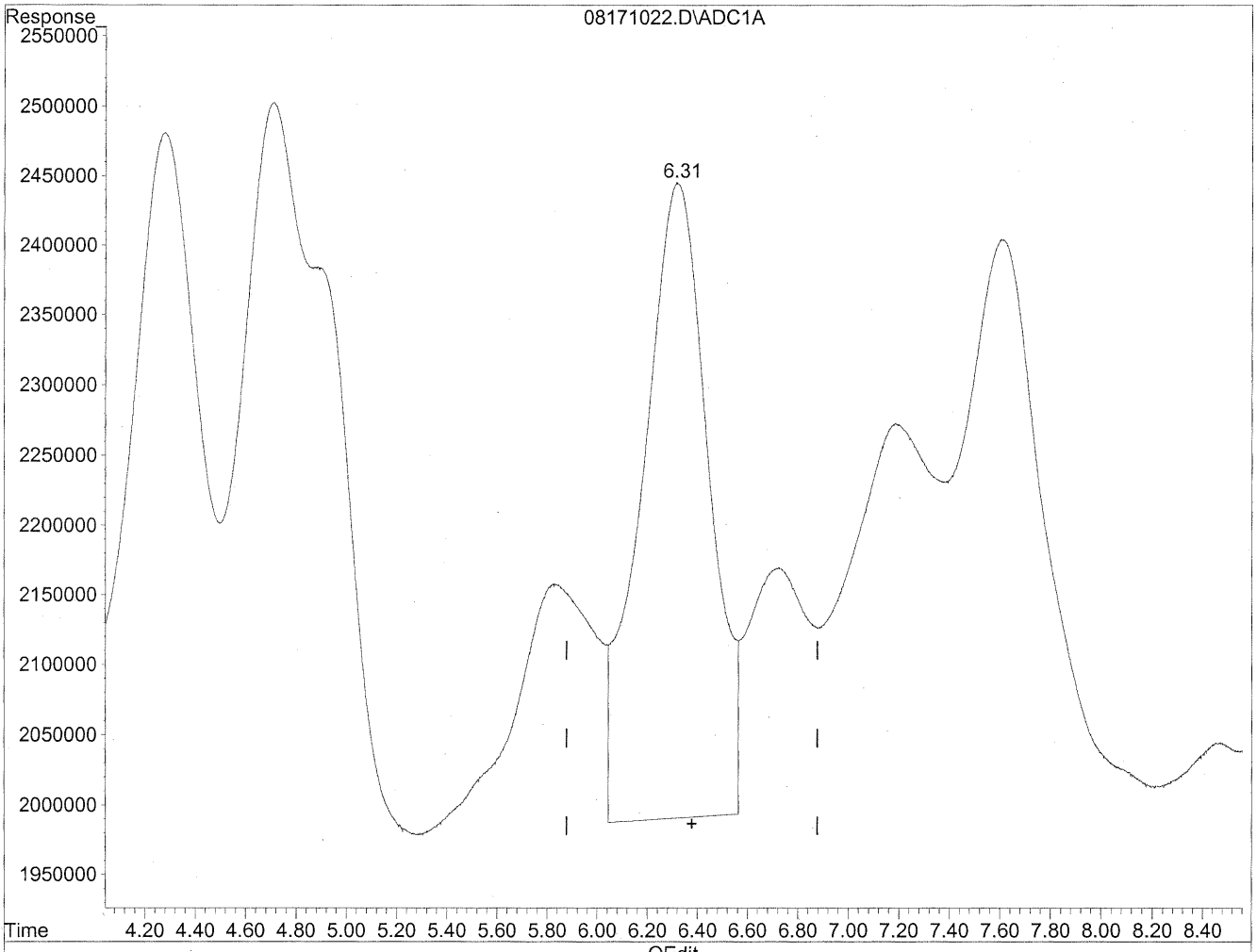
296

(+) = Expected Retention Time
08171022.D TO110709.M Sat Aug 22 15:30:08 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:28 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(6) Benzaldehyde

6.32min 1278.310ng/ml

response 84201373

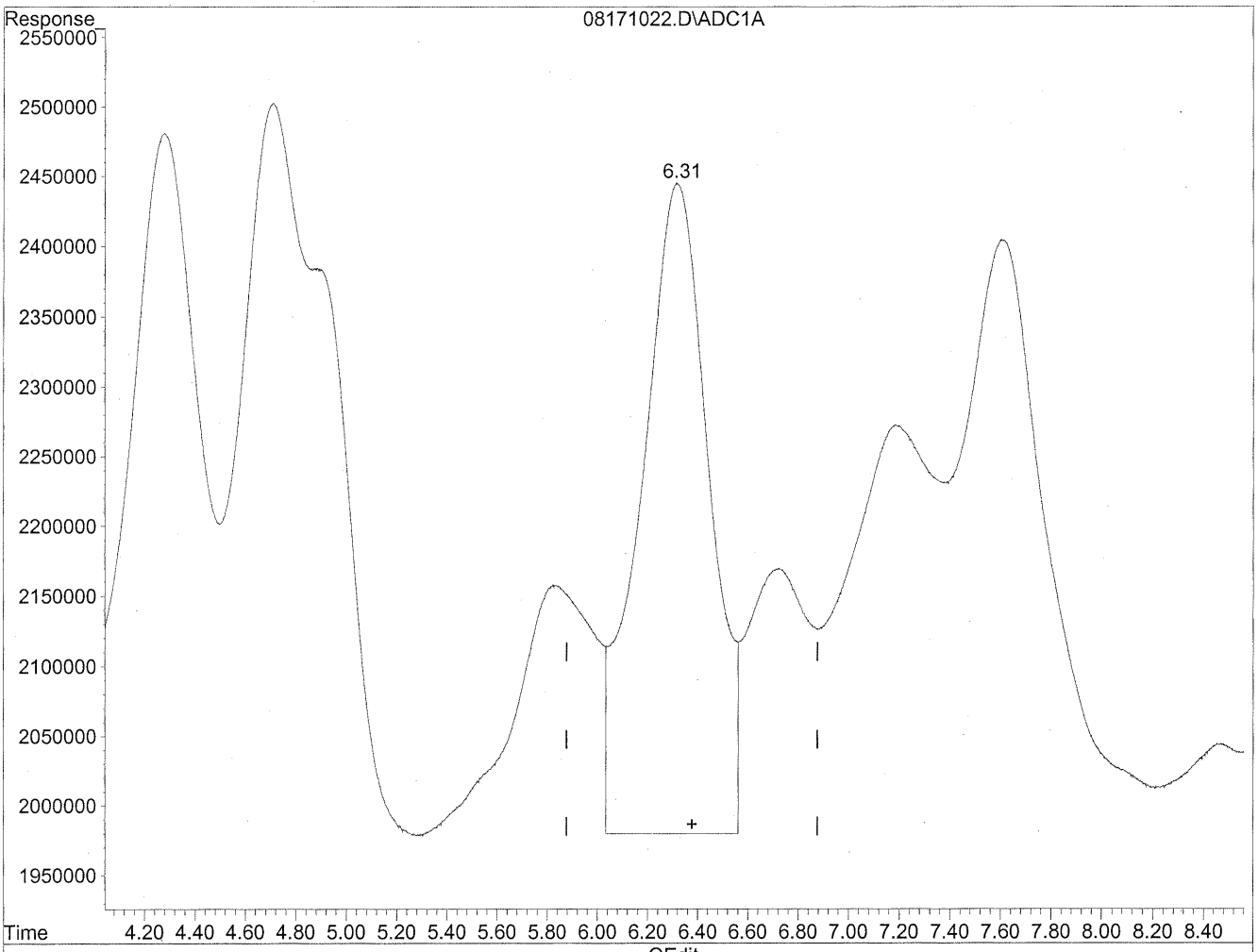
297

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:28 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.31min 1337.327ng/ml m

response 88088777

HC 8/22/09
BC

KF 8/23/09

298

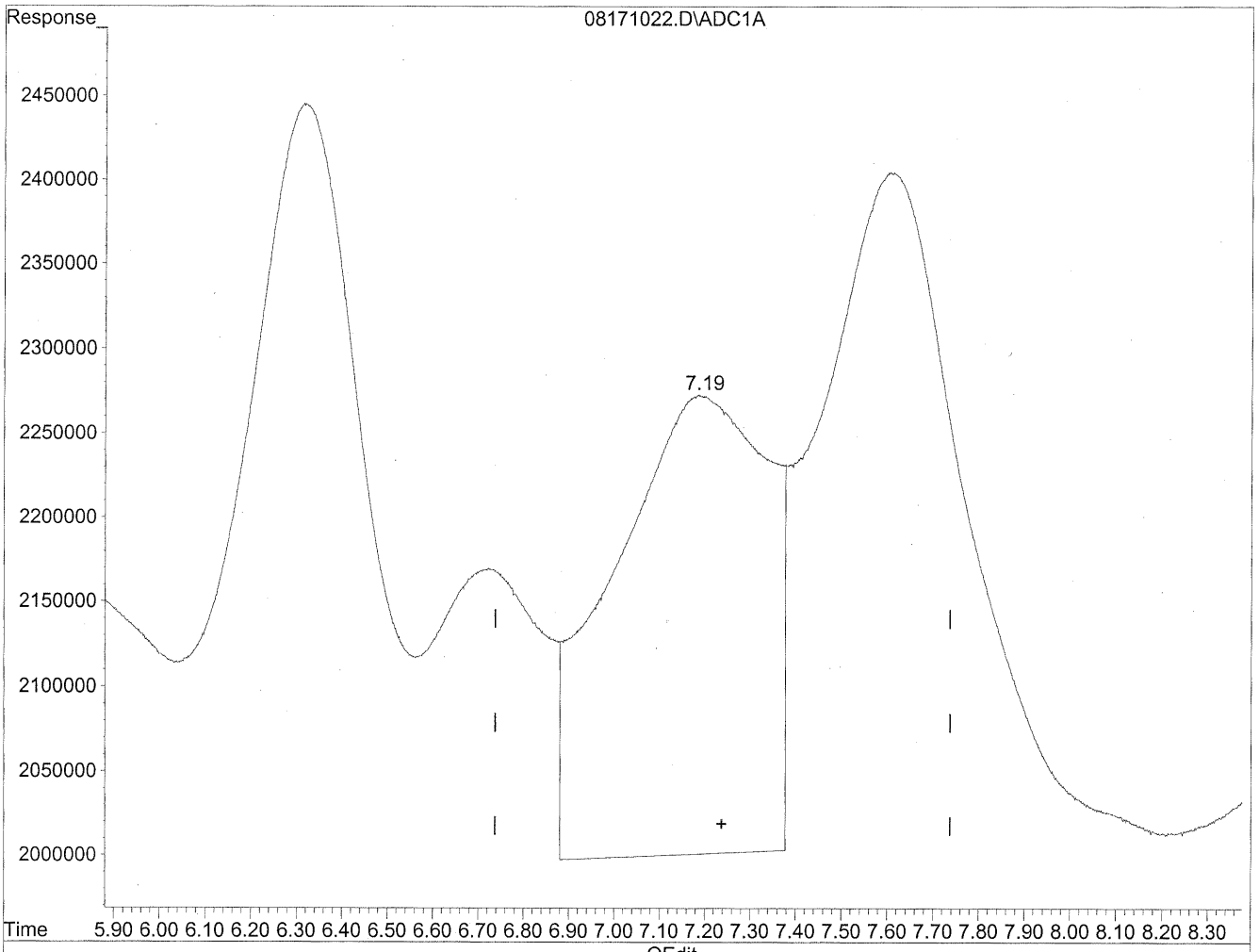
(+) = Expected Retention Time

08171022.D TO110709.M Sat Aug 22 15:30:57 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:28 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.19min 821.355ng/ml

response 64271854

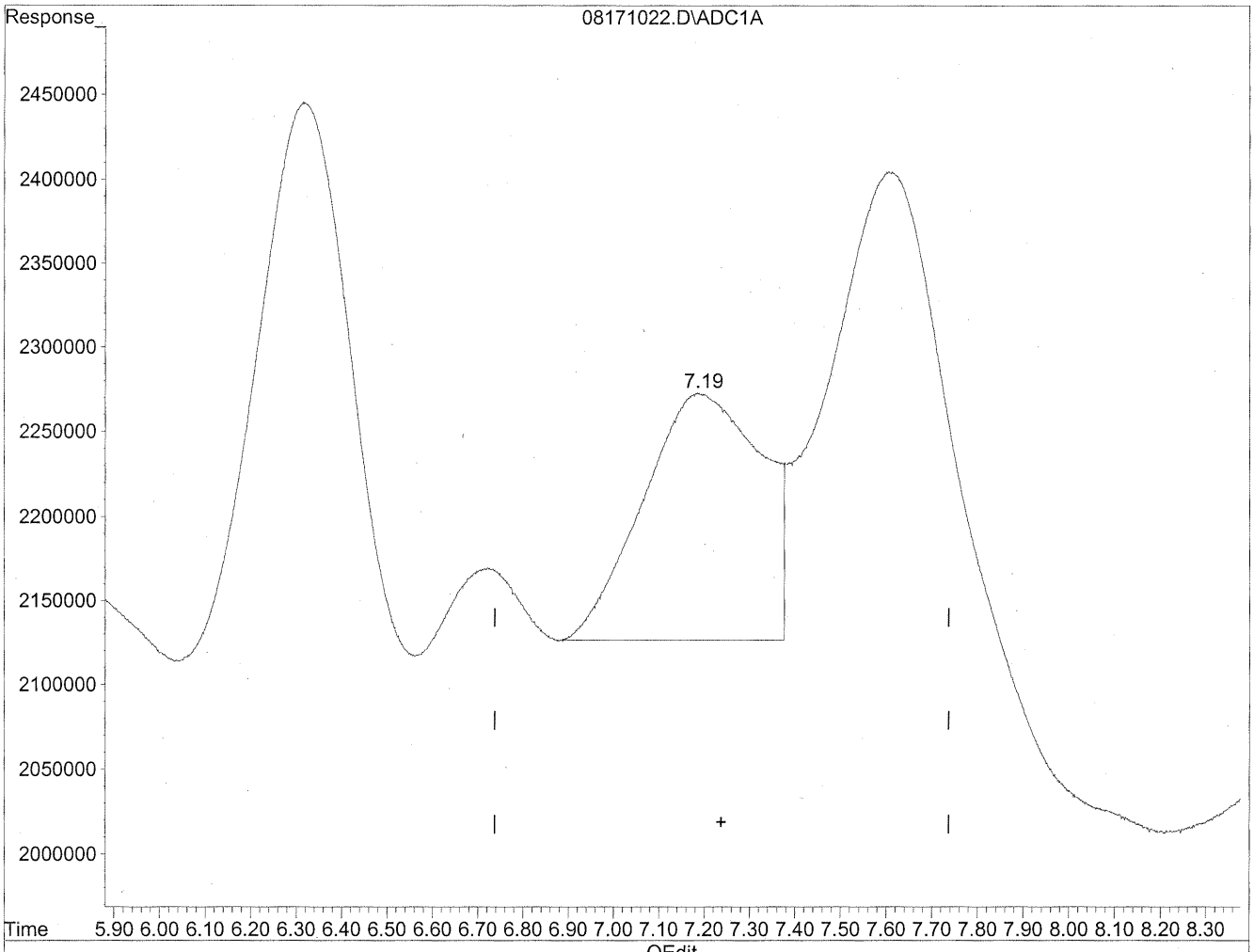
299

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:28 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.19min 339.469ng/ml m

response 26563770

HC 8/22/09
BC

KK 8/23/09

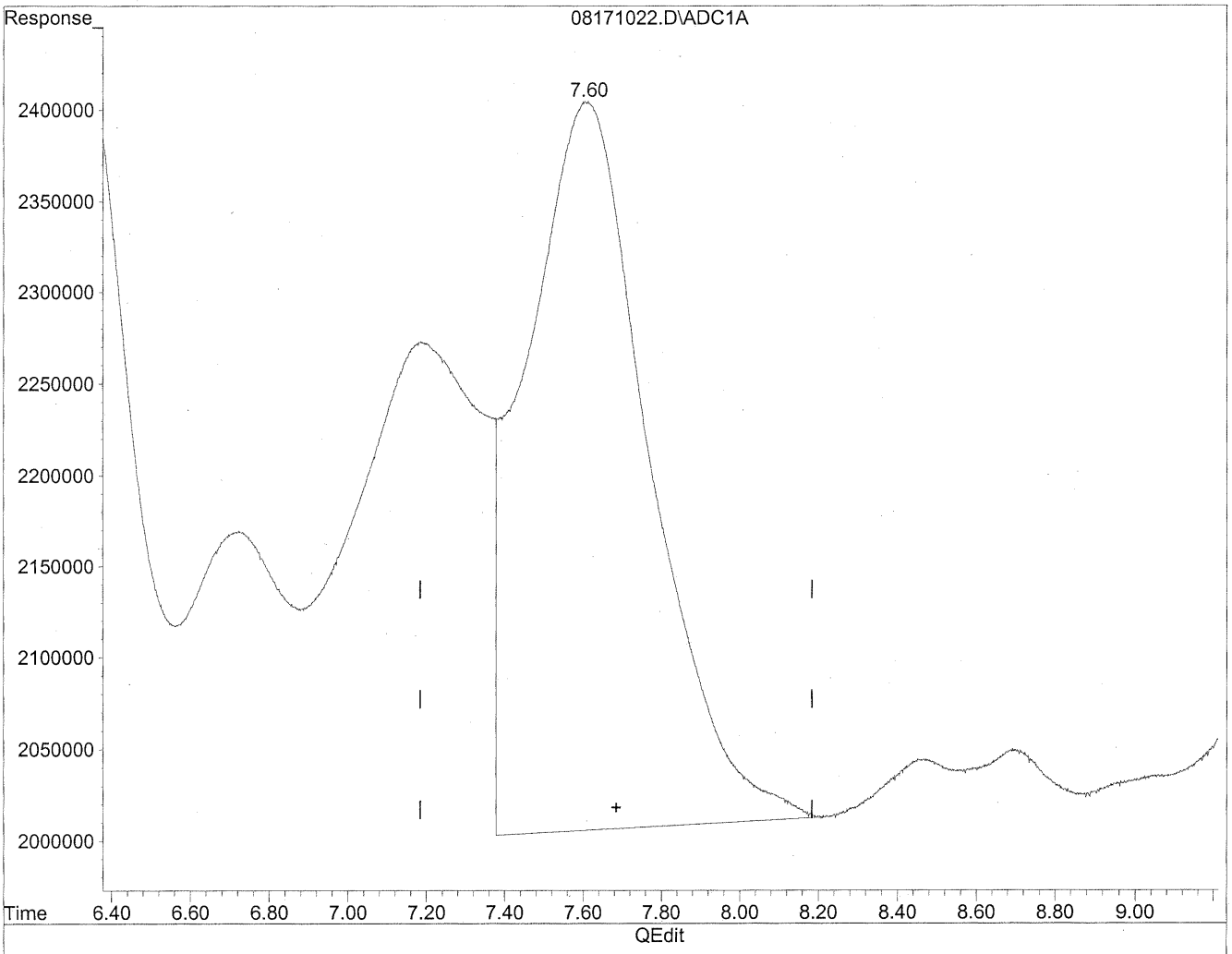
(+) = Expected Retention Time

08171022.D TO110709.M Sat Aug 22 15:31:29 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:28 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.61min 1192.452ng/ml

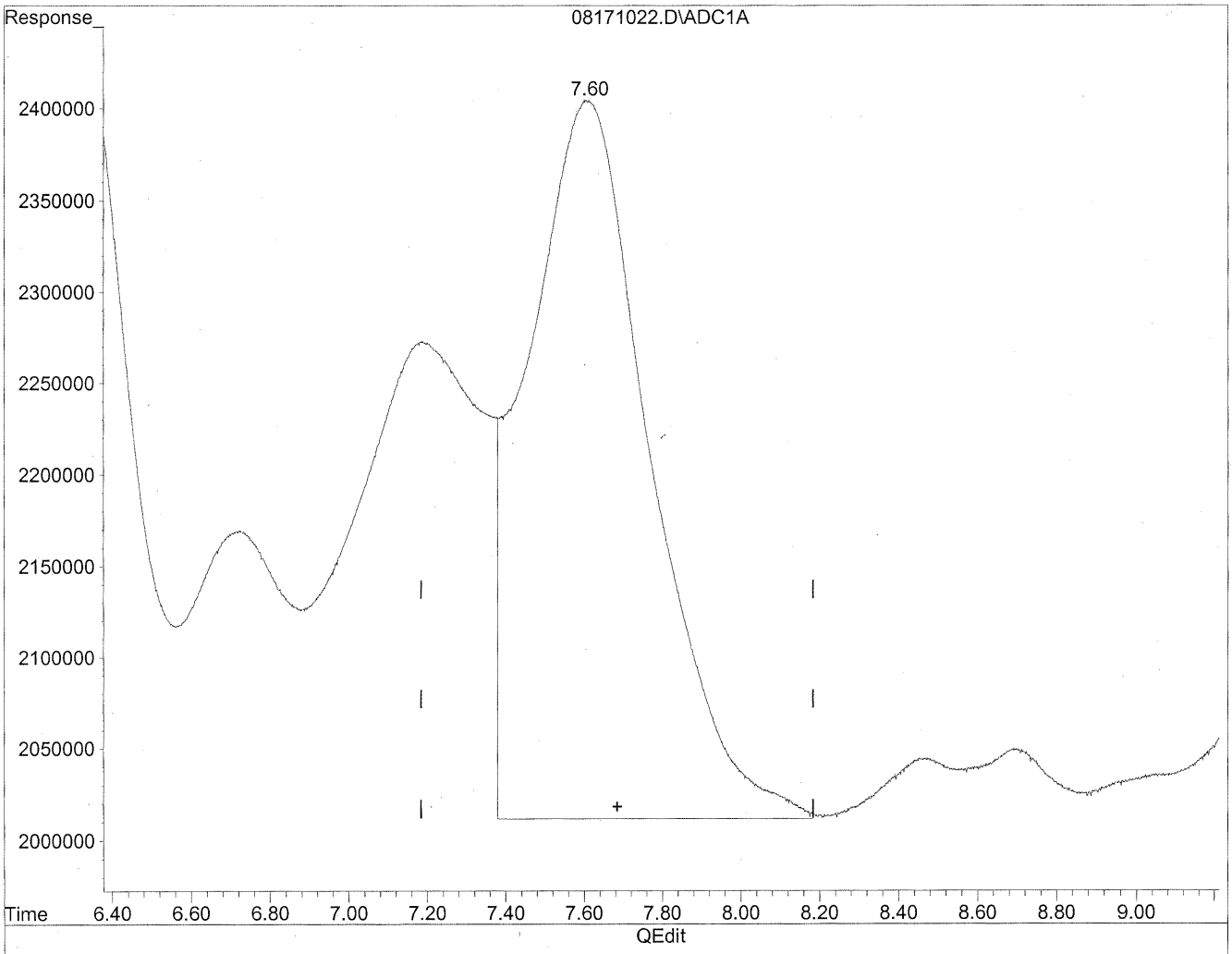
response 87651175

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:28 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.60min 1163.409ng/ml m

response 85516374

HC 8/22/09 BC

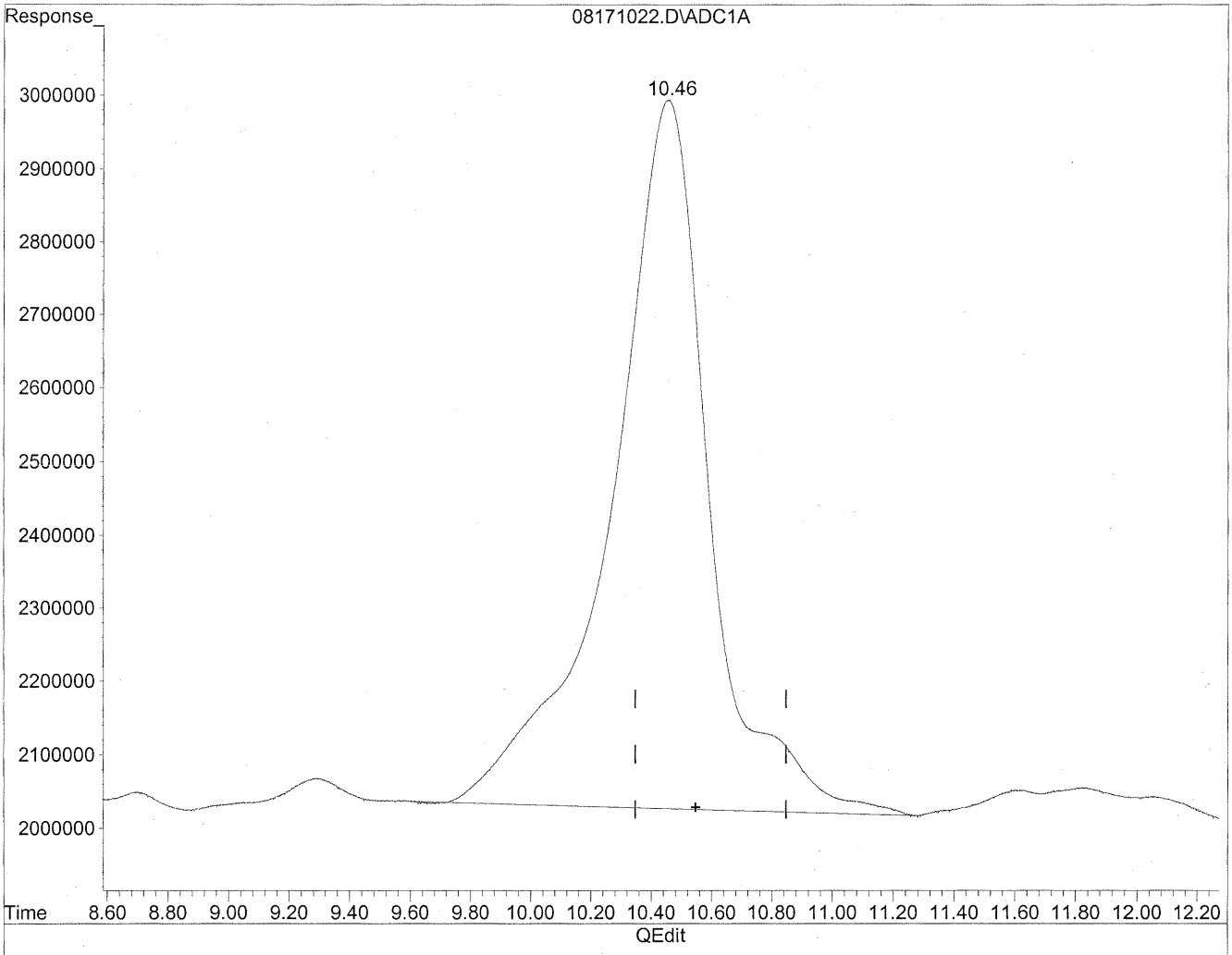
KR 8/23/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:28 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.46min 3211.797ng/ml

response 216294514

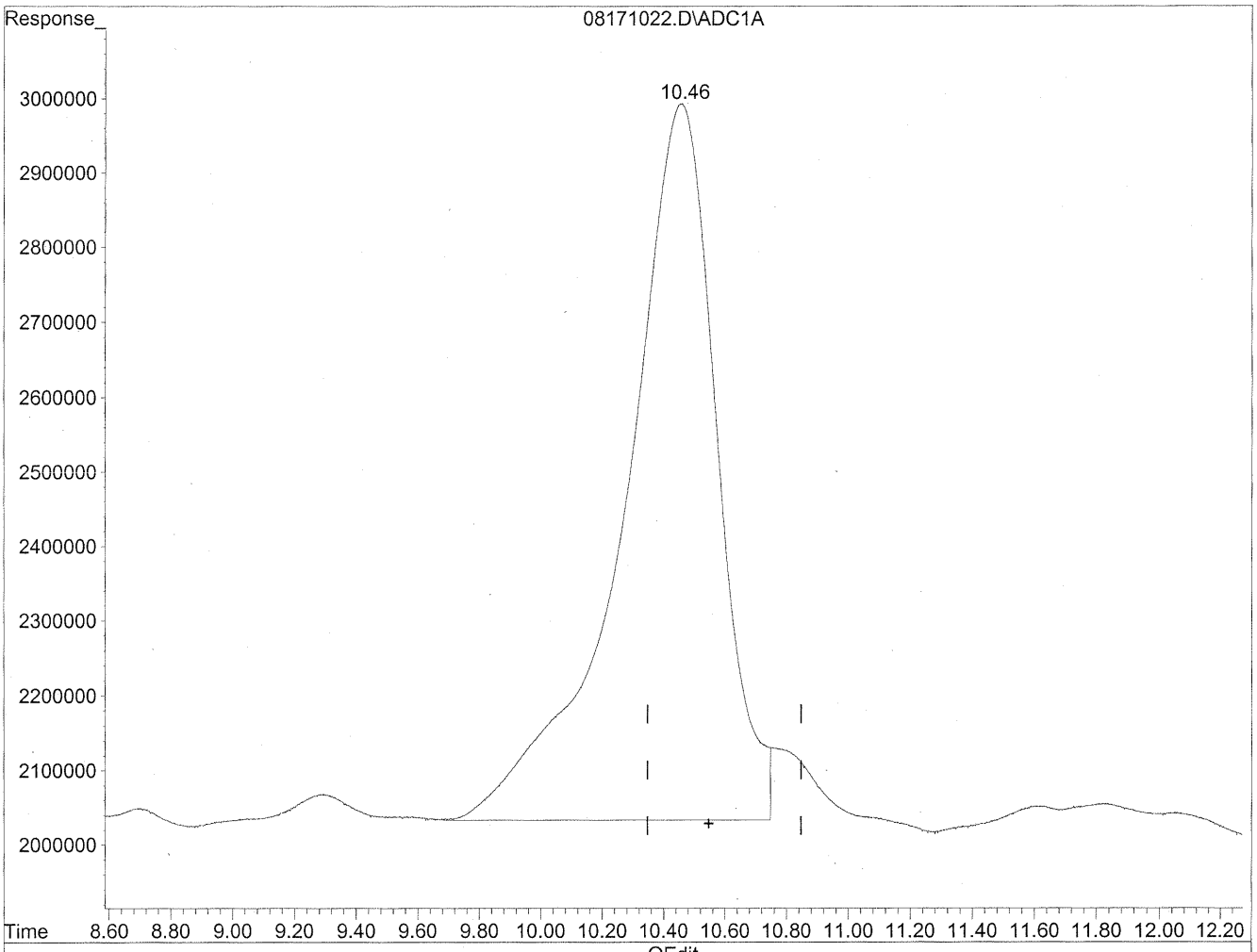
303

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:28 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.46min 2989.804ng/ml m

response 201344667

tlc 8/22/09 SH

KR 8/23/09

304

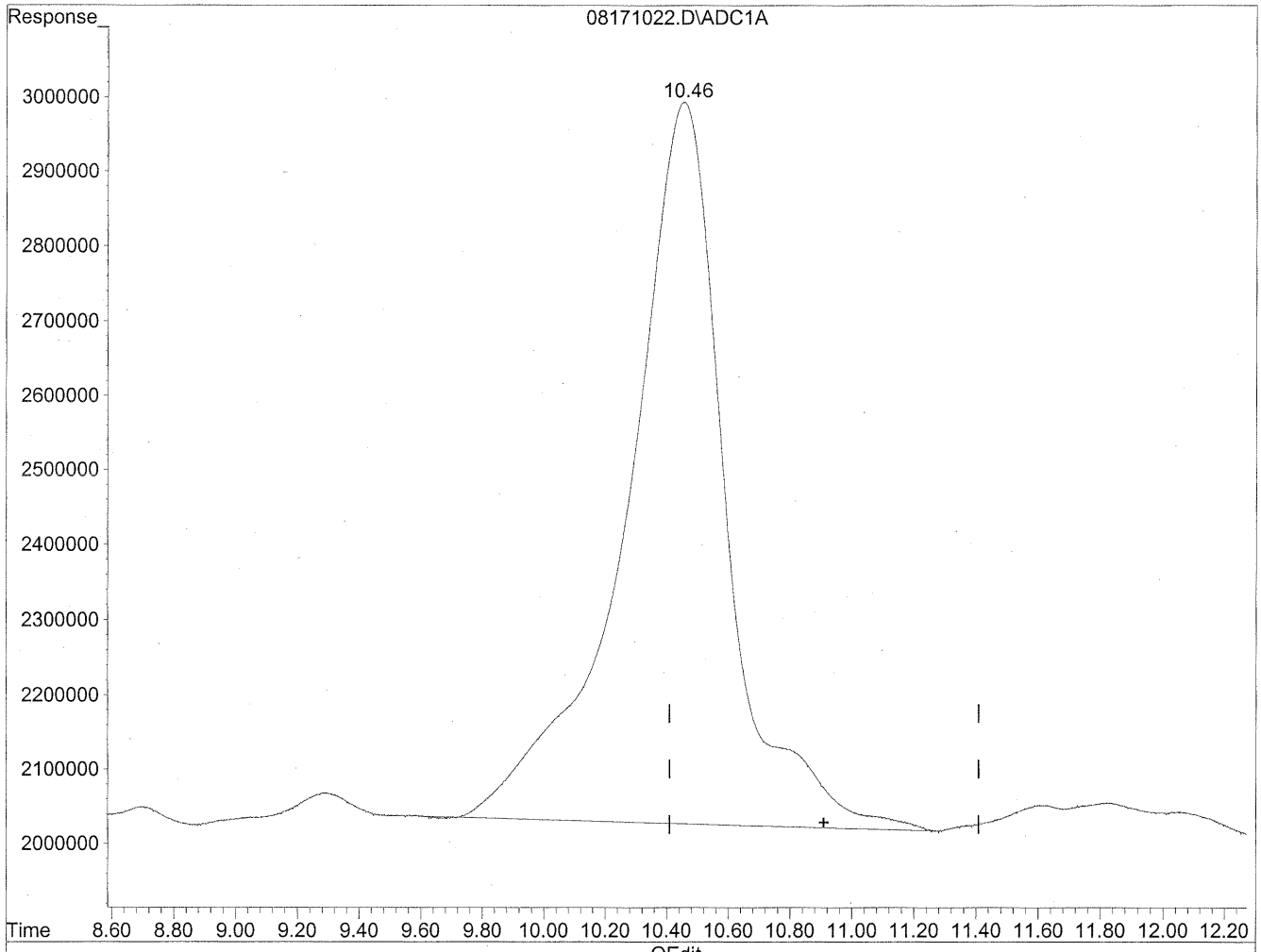
(+) = Expected Retention Time

08171022.D TO110709.M Sat Aug 22 15:32:27 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:28 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.46min 4412.968ng/ml

response 216294514

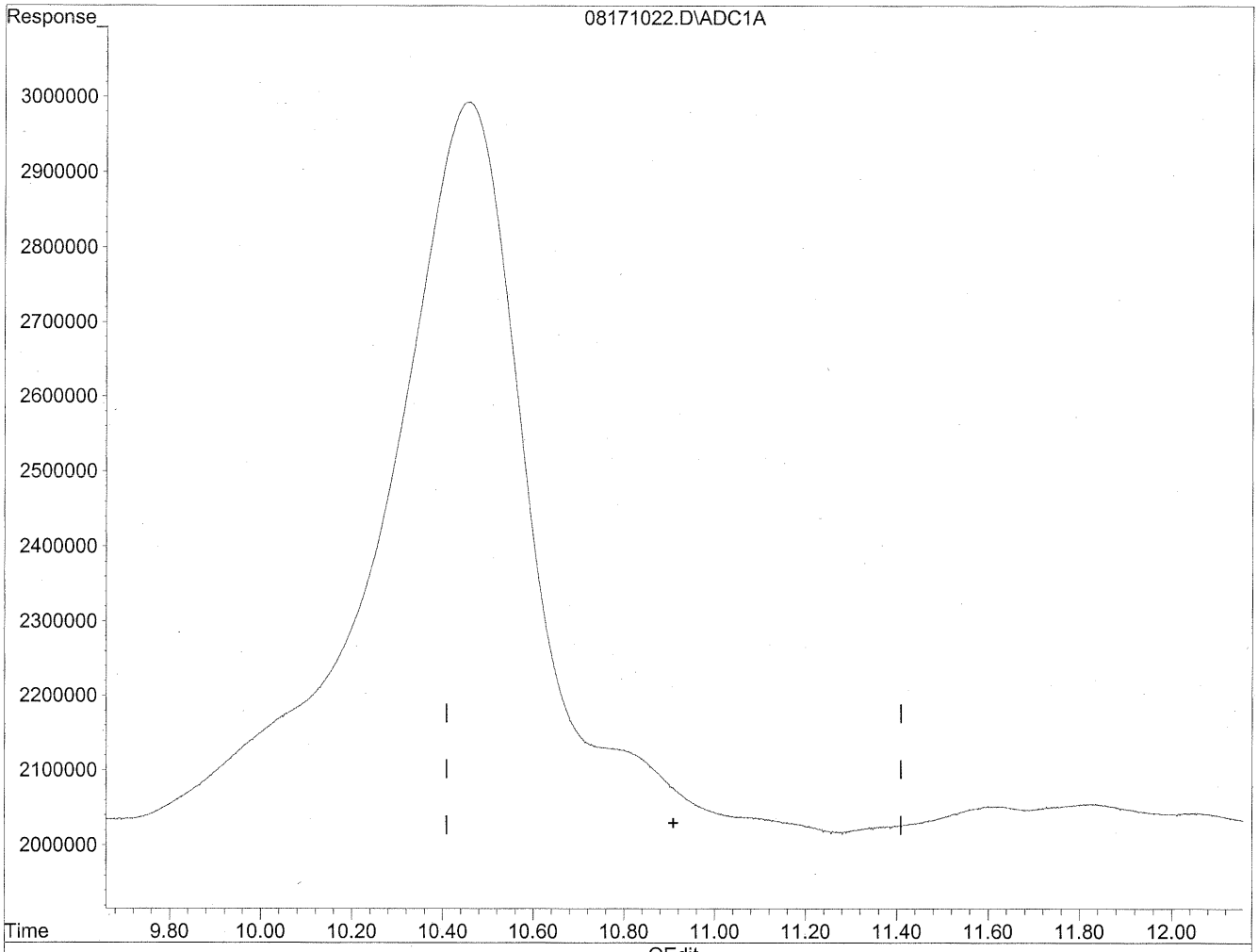
(+) = Expected Retention Time

08171022.D TO110709.M Sat Aug 22 15:32:32 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171022.D Vial: 35
Acq On : 18 Aug 2009 9:09 pm Operator: HC
Sample : P0902772-012 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 15:28 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/22/09
MP
KP 8/23/09

306

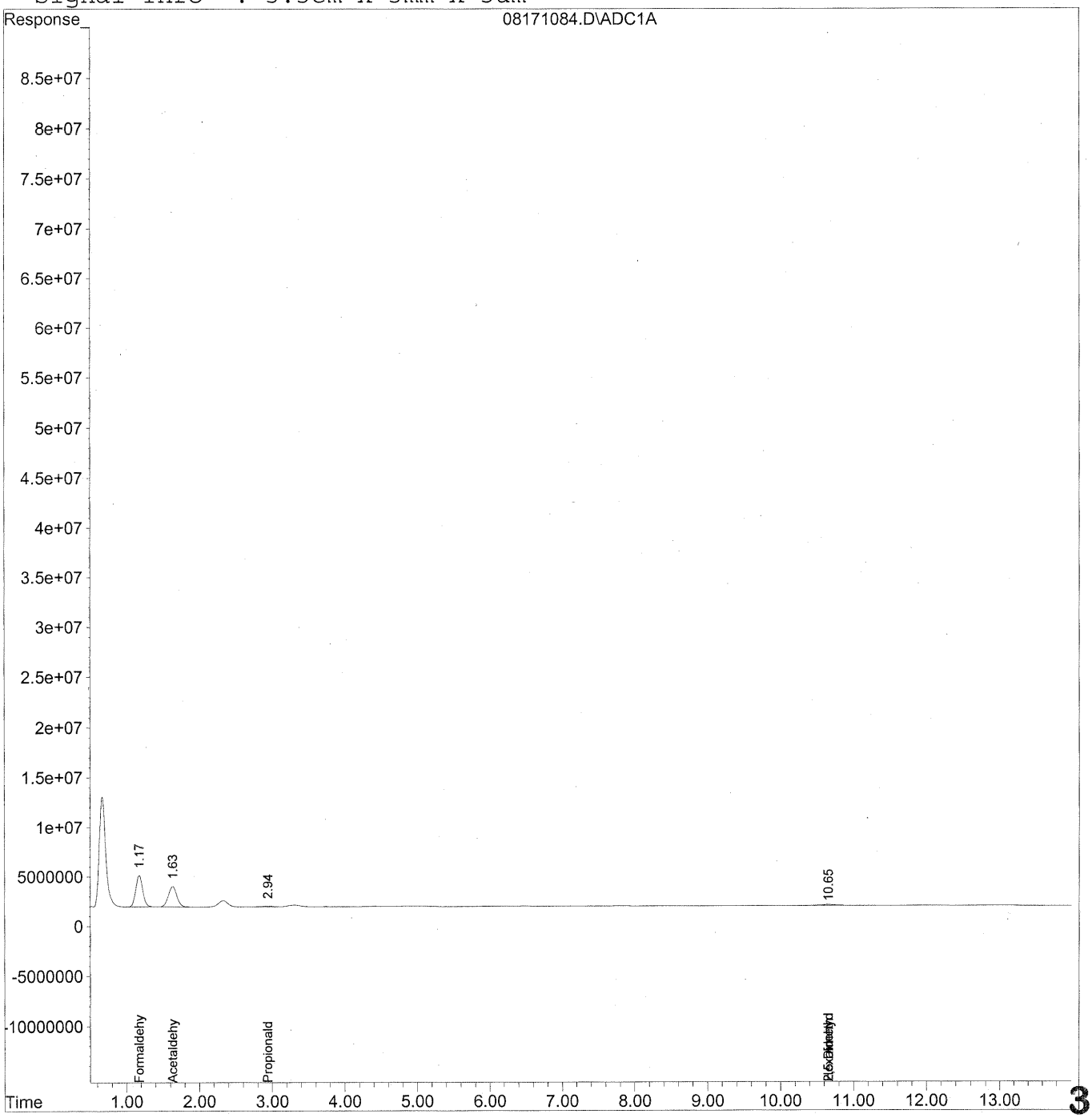
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171084.D Vial: 9
Acq On : 19 Aug 2009 12:42 pm Operator: HC
Sample : P0902772-012 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:23 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



307

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171084.D Vial: 9
Acq On : 19 Aug 2009 12:42 pm Operator: HC
Sample : P0902772-012 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:23 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

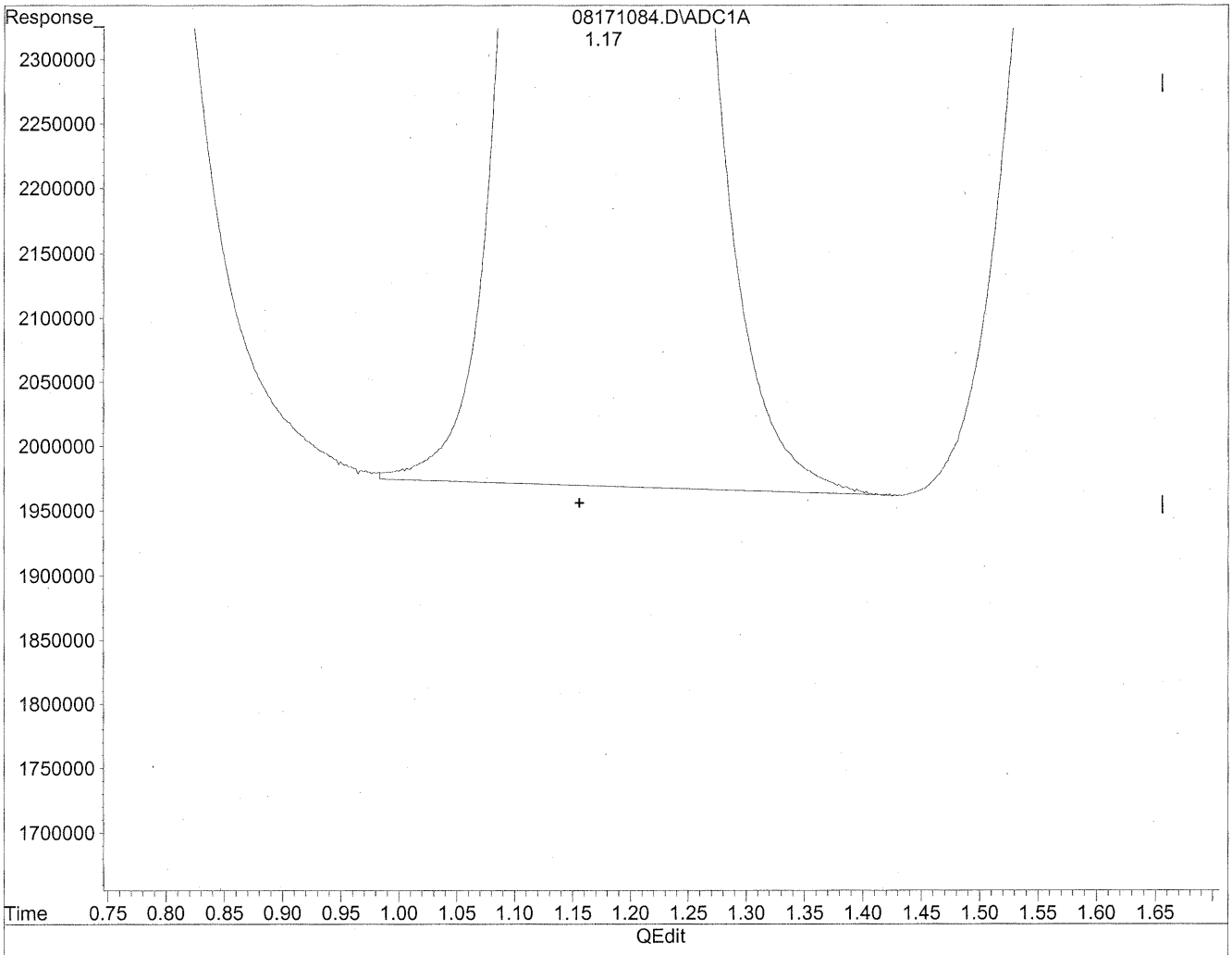
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	211781944	1153.613	ng/mlm
2)	Acetaldehyde	1.63	169763661	1210.665	ng/ml
3)	Propionaldehyde	2.94	8740850	81.924	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	10.66	17805860	264.403	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.66f	17805860	363.286	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171084.D Vial: 9
Acq On : 19 Aug 2009 12:42 pm Operator: HC
Sample : P0902772-012 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 1156.185ng/ml

response 212254120

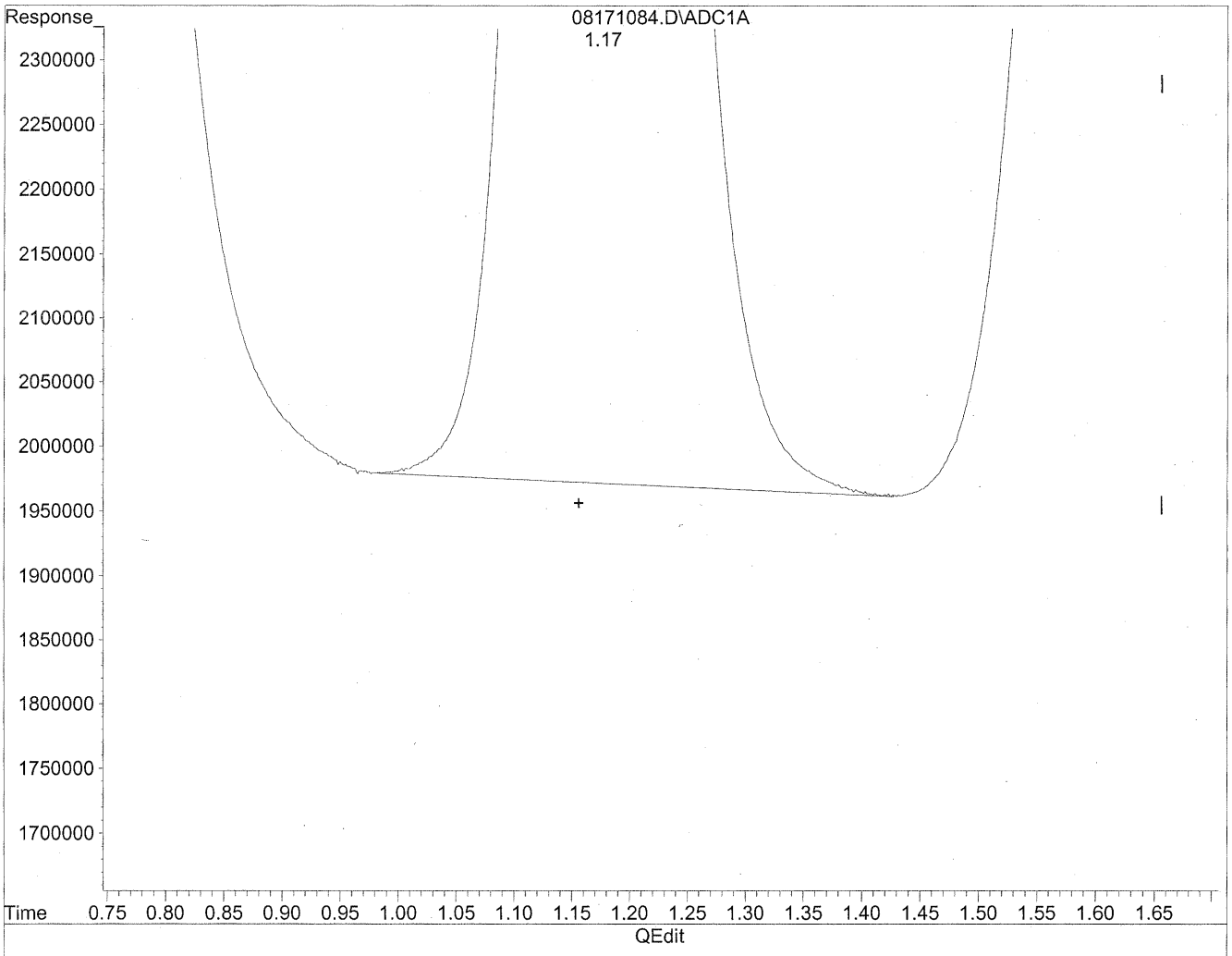
(+) = Expected Retention Time

08171084.D TO110709.M Sat Aug 22 17:23:42 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171084.D Vial: 9
Acq On : 19 Aug 2009 12:42 pm Operator: HC
Sample : P0902772-012 front 10x Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 17:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 1153.613ng/ml m

response 211781944

HC 8/22/09 LC

KR 8/23/09

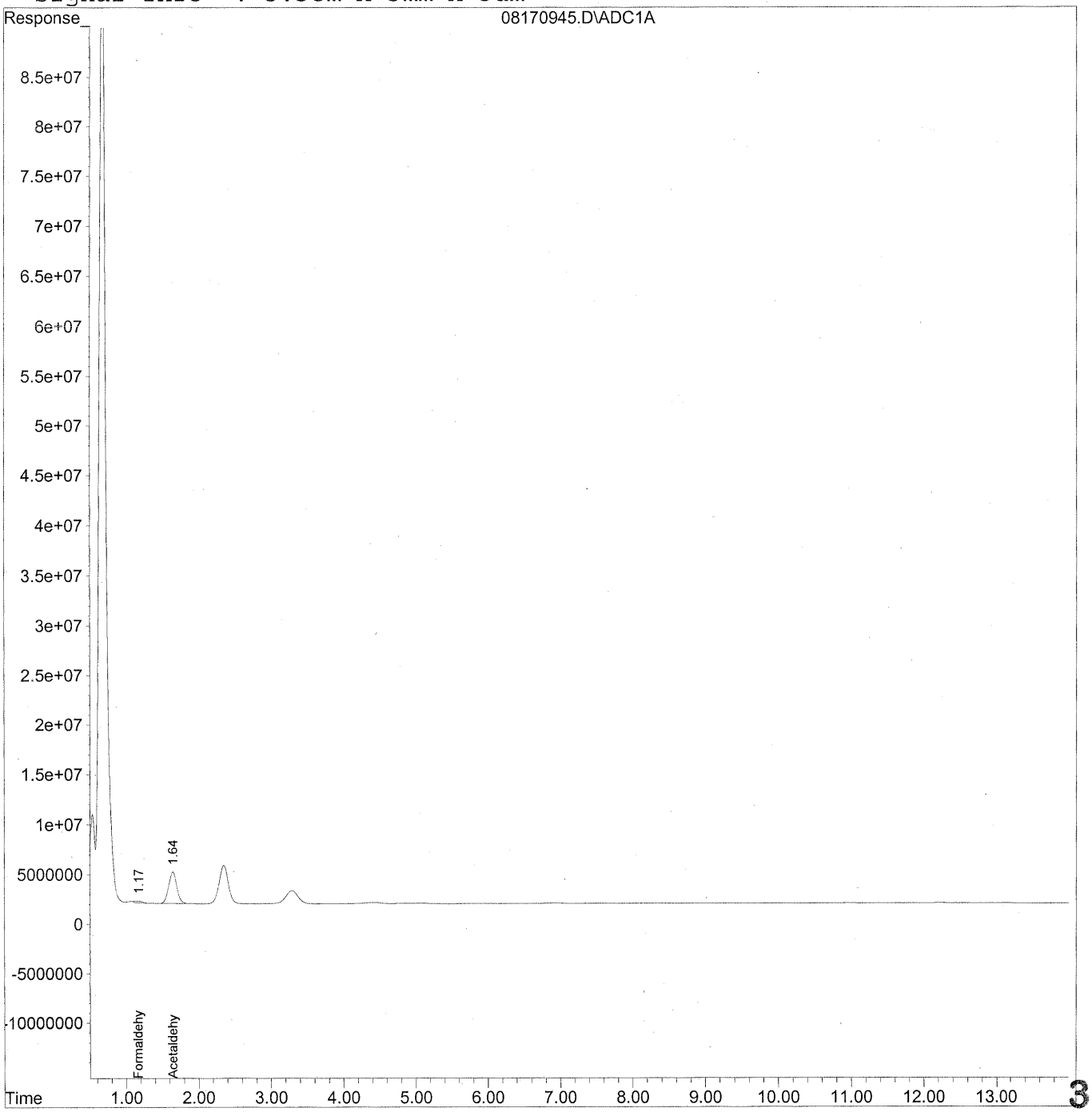
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170945.D Vial: 44
Acq On : 18 Aug 2009 1:52 am Operator: HC
Sample : P0902772-012 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:36 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



311

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170945.D Vial: 44
Acq On : 18 Aug 2009 1:52 am Operator: HC
Sample : P0902772-012 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:36 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

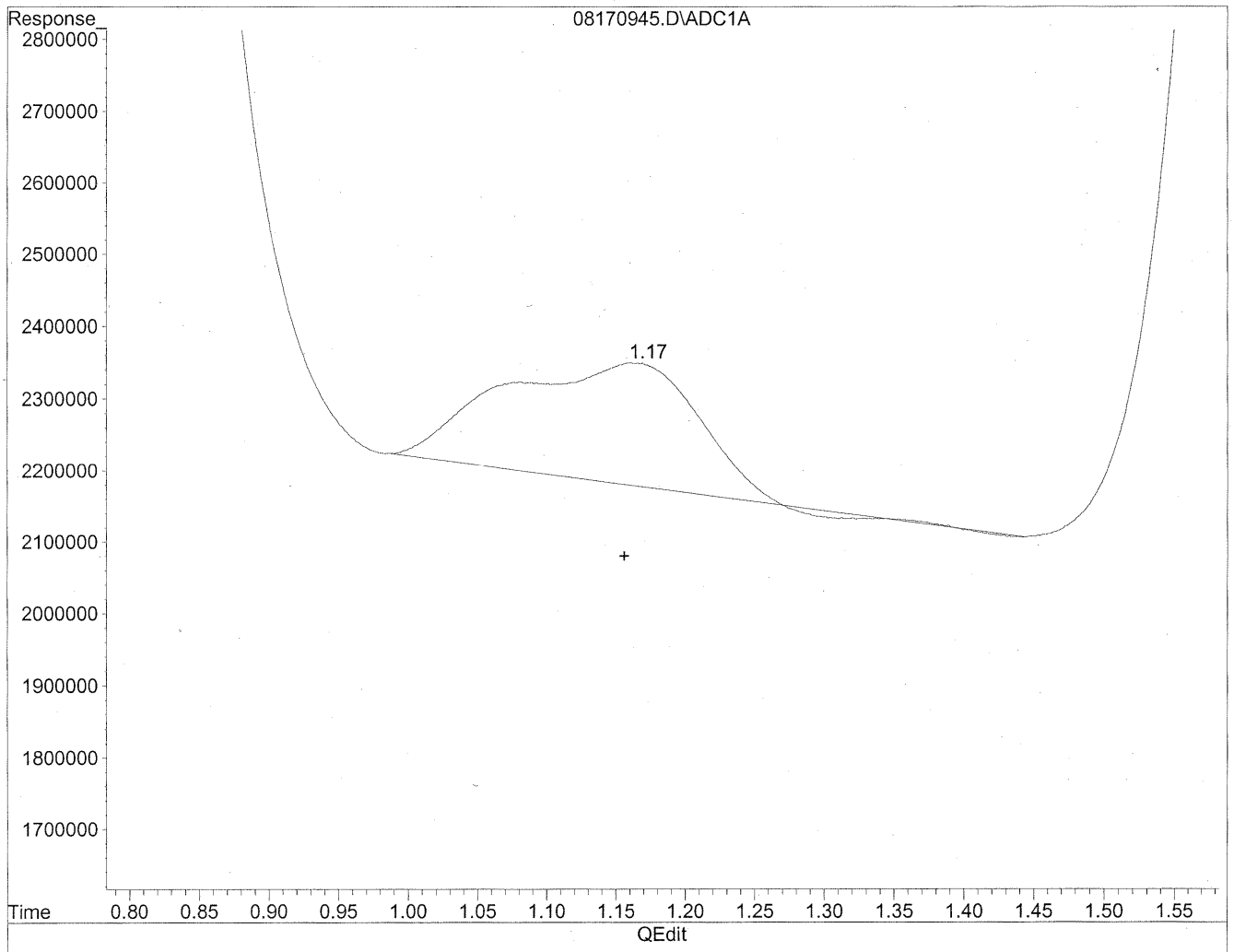
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	11638436	63.397	ng/mlm
2)	Acetaldehyde	1.64	251737990	1795.262	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170945.D Vial: 44
Acq On : 18 Aug 2009 1:52 am Operator: HC
Sample : P0902772-012 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:36 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.16min 87.952ng/ml

response 16146424

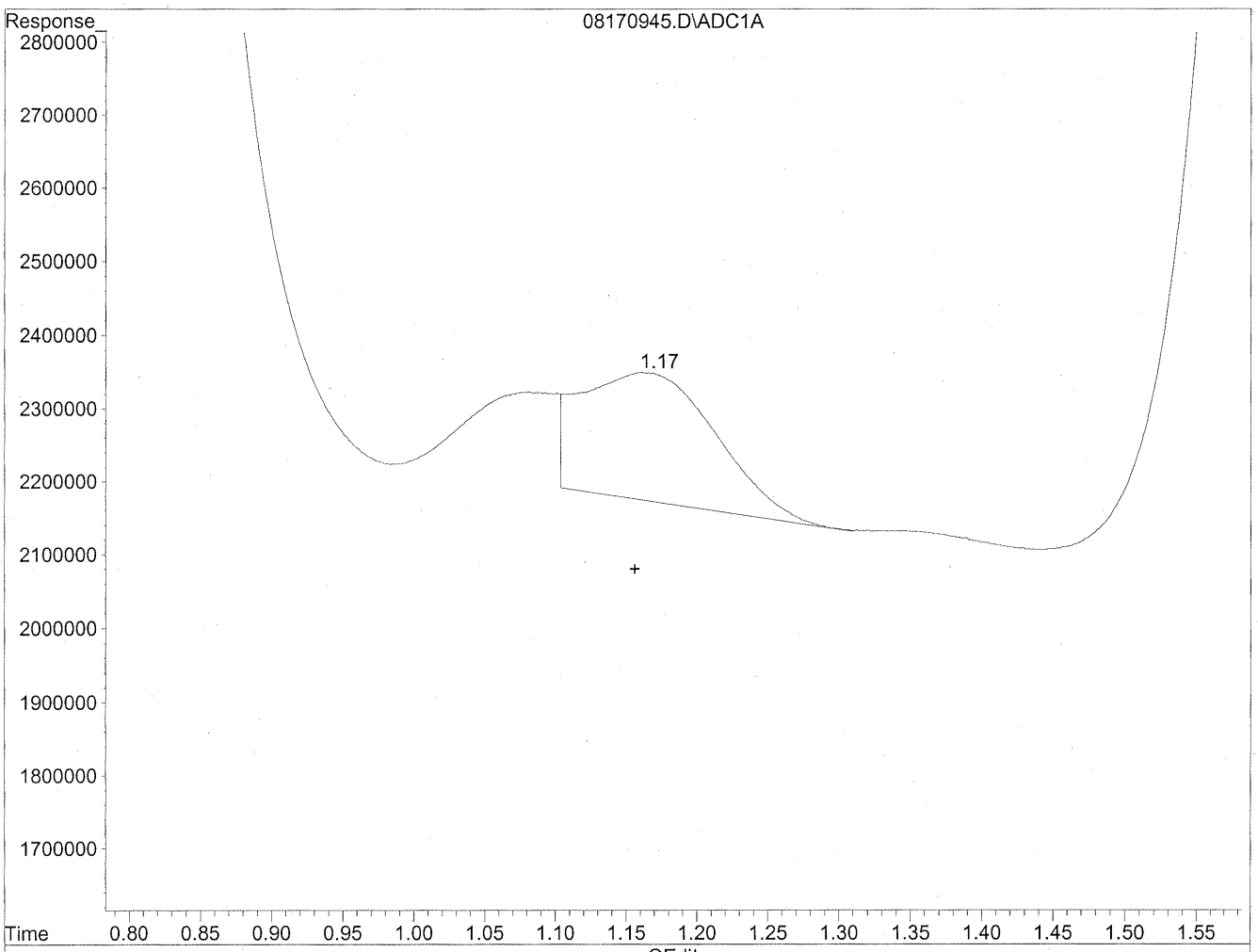
313

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170945.D Vial: 44
Acq On : 18 Aug 2009 1:52 am Operator: HC
Sample : P0902772-012 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:36 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 63.397ng/ml m

response 11638436

HC 8/22/09
SP

KE 8/23/09

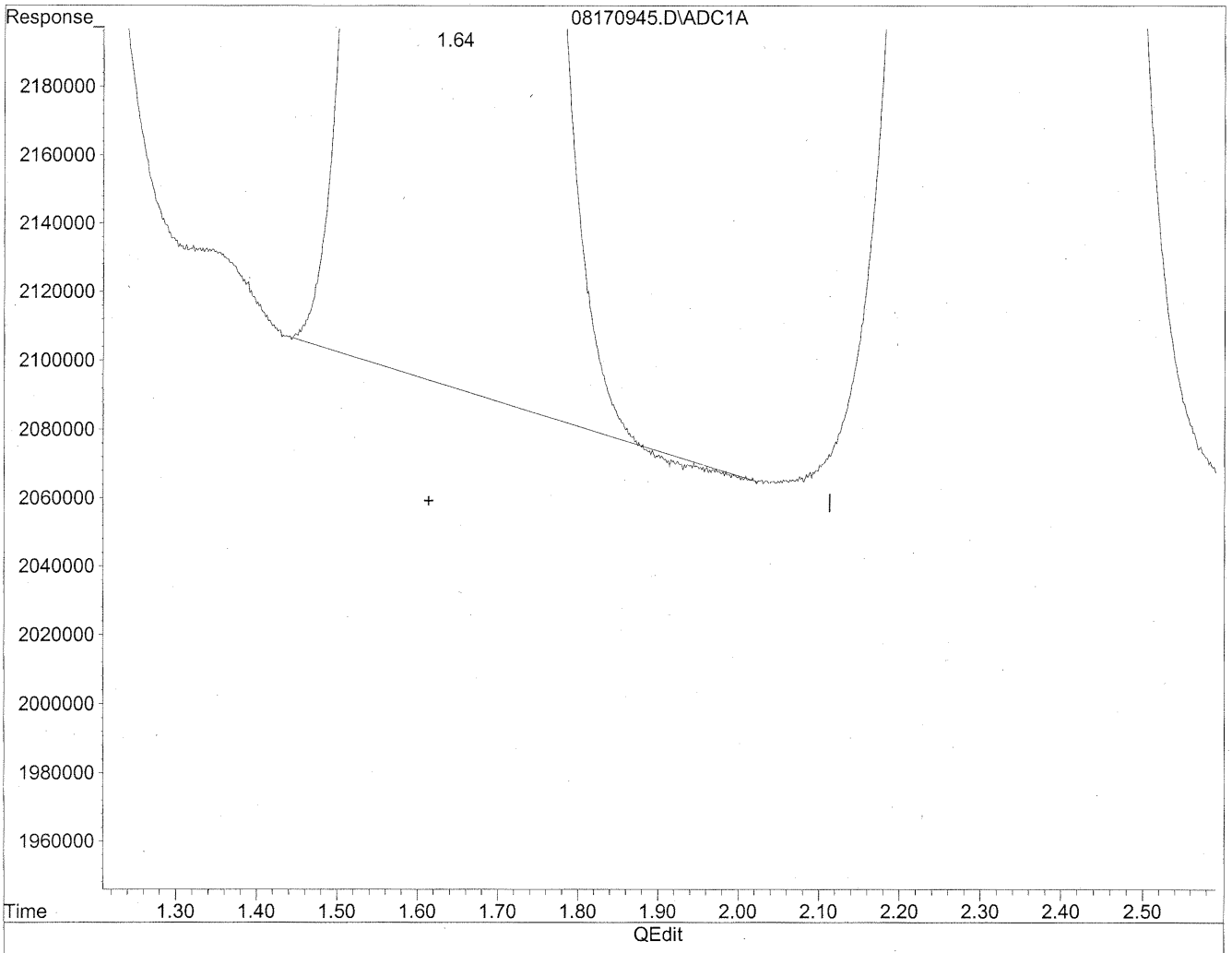
314

(+) = Expected Retention Time
08170945.D TO110709.M Sat Aug 22 12:36:27 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170945.D Vial: 44
Acq On : 18 Aug 2009 1:52 am Operator: HC
Sample : P0902772-012 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:36 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.64min 1791.891ng/ml

response 251265227

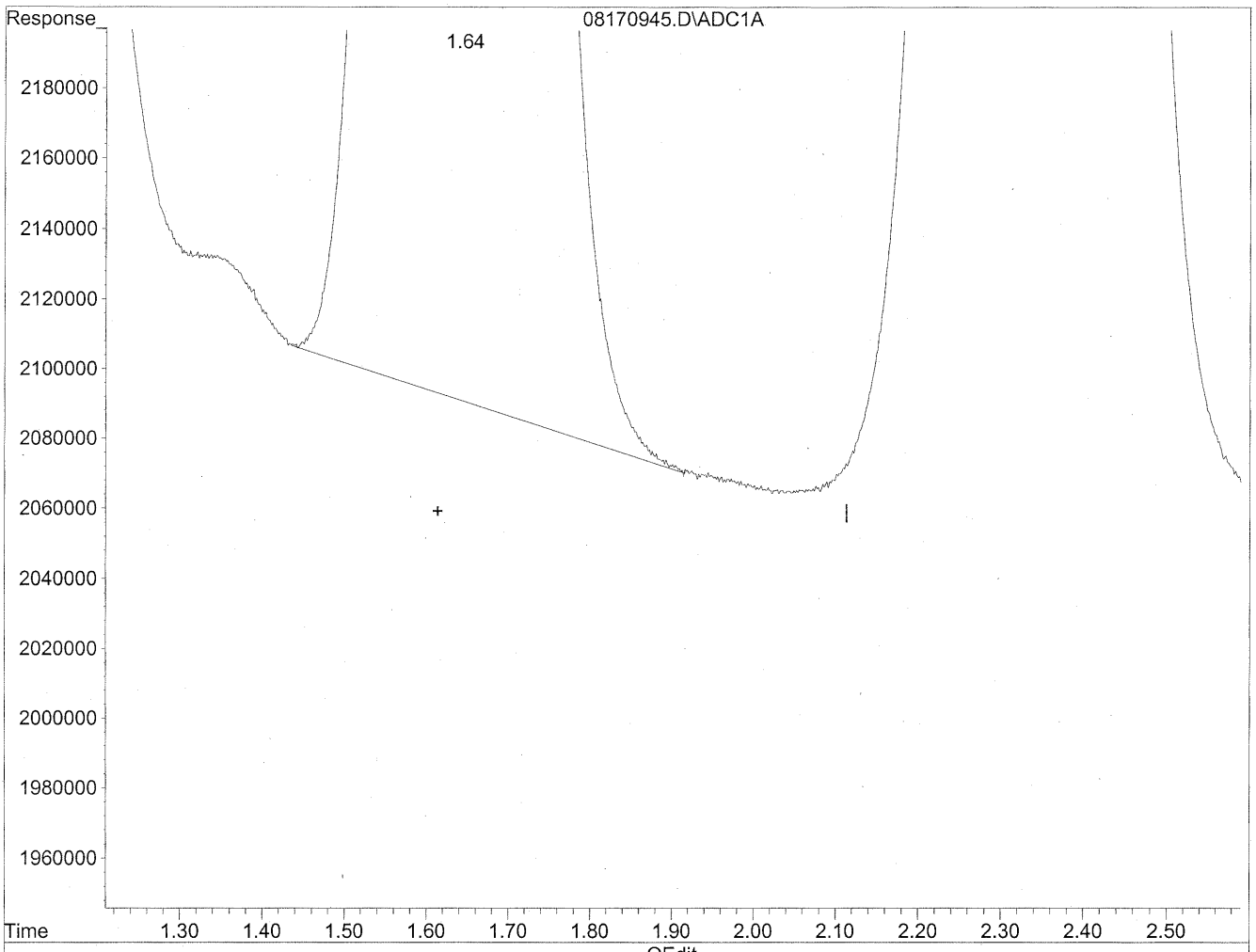
315

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170945.D Vial: 44
Acq On : 18 Aug 2009 1:52 am Operator: HC
Sample : P0902772-012 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:36 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.64min 1795.262ng/ml m

response 251737990

HC 8/24/09
LC
pp 8/23/09

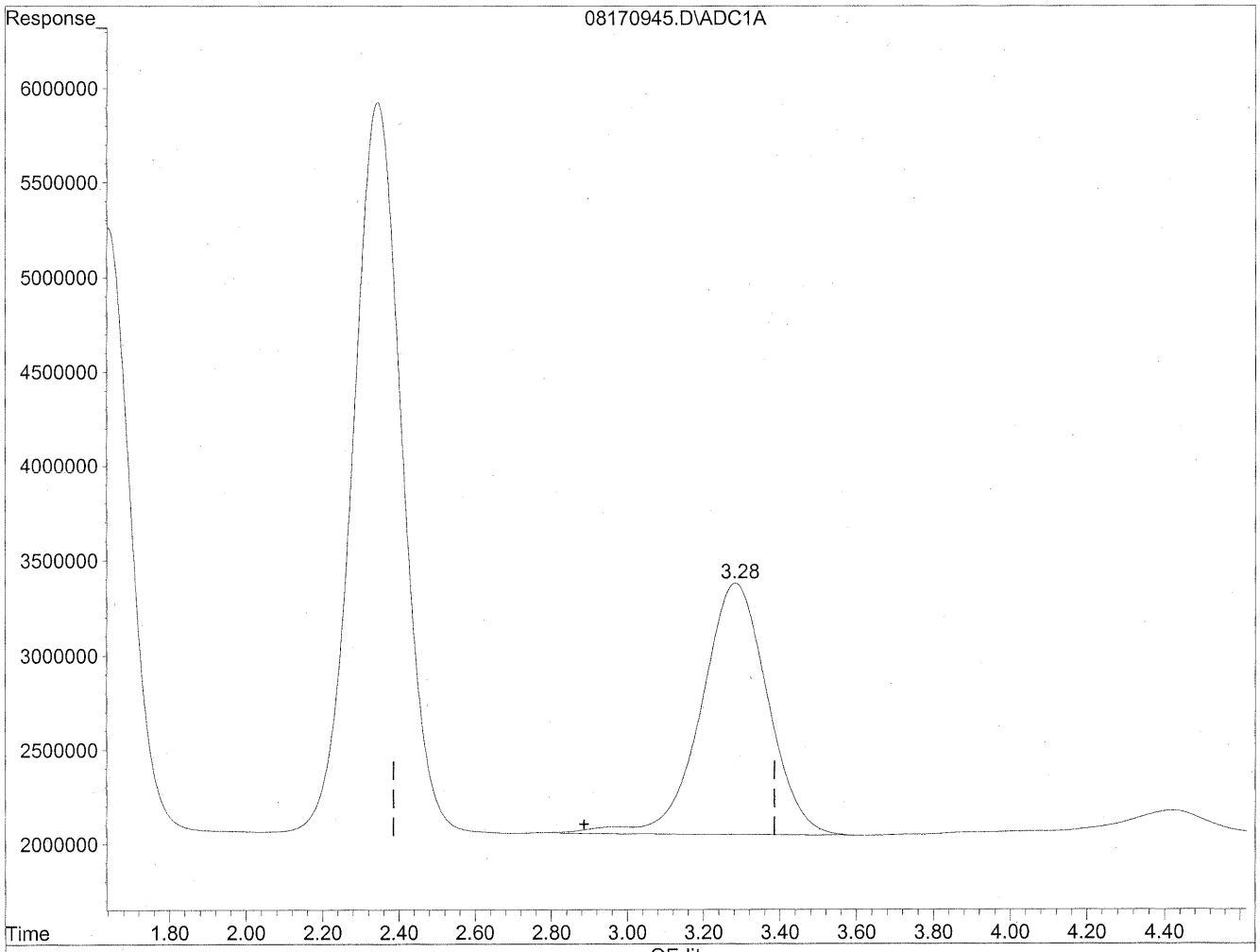
316

(+) = Expected Retention Time
08170945.D TO110709.M Sat Aug 22 12:36:44 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170945.D Vial: 44
Acq On : 18 Aug 2009 1:52 am Operator: HC
Sample : P0902772-012 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:36 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.28min 1508.910ng/ml

response 160993463

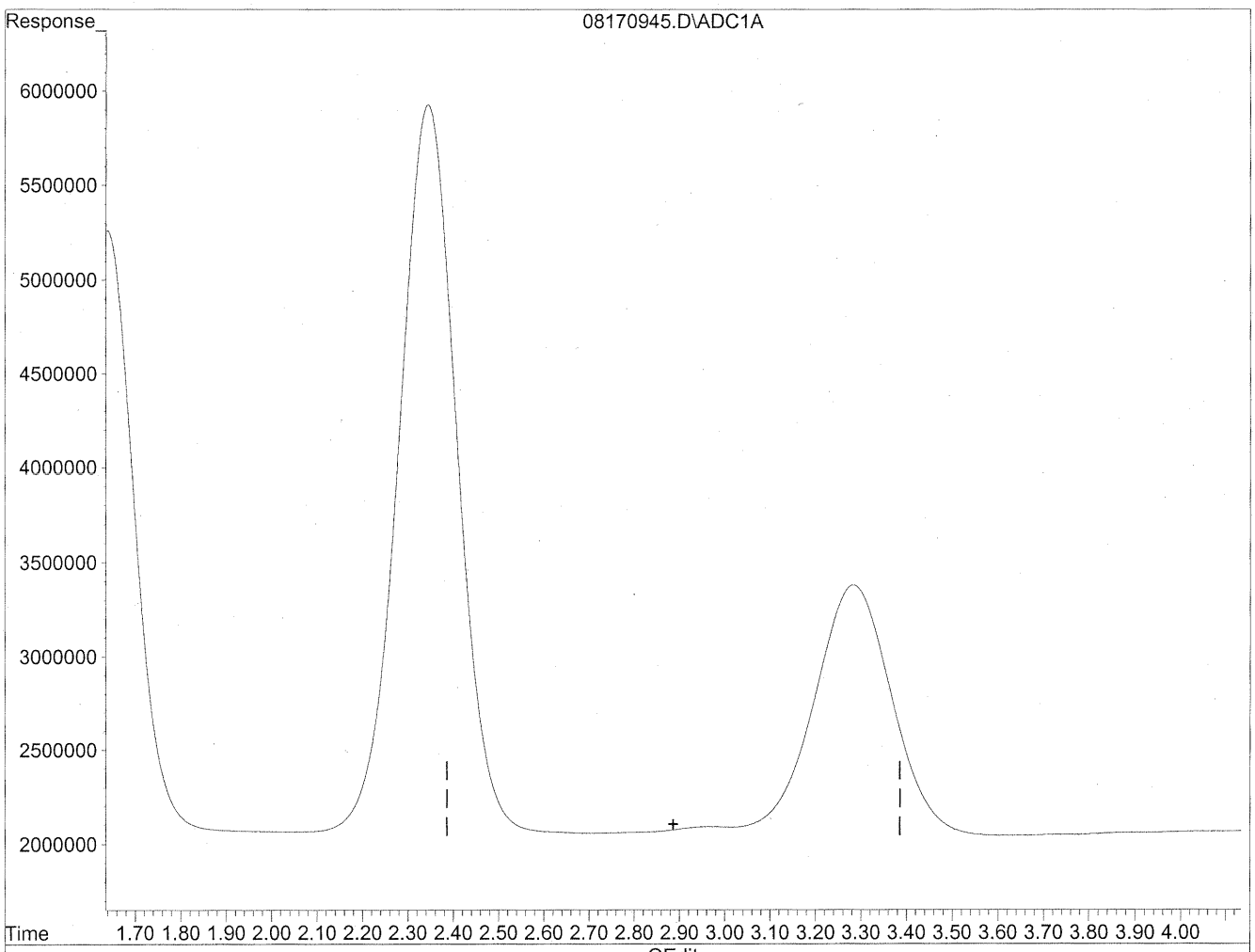
317

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170945.D Vial: 44
Acq On : 18 Aug 2009 1:52 am Operator: HC
Sample : P0902772-012 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:36 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/24/09 MP

KK 8/23/09

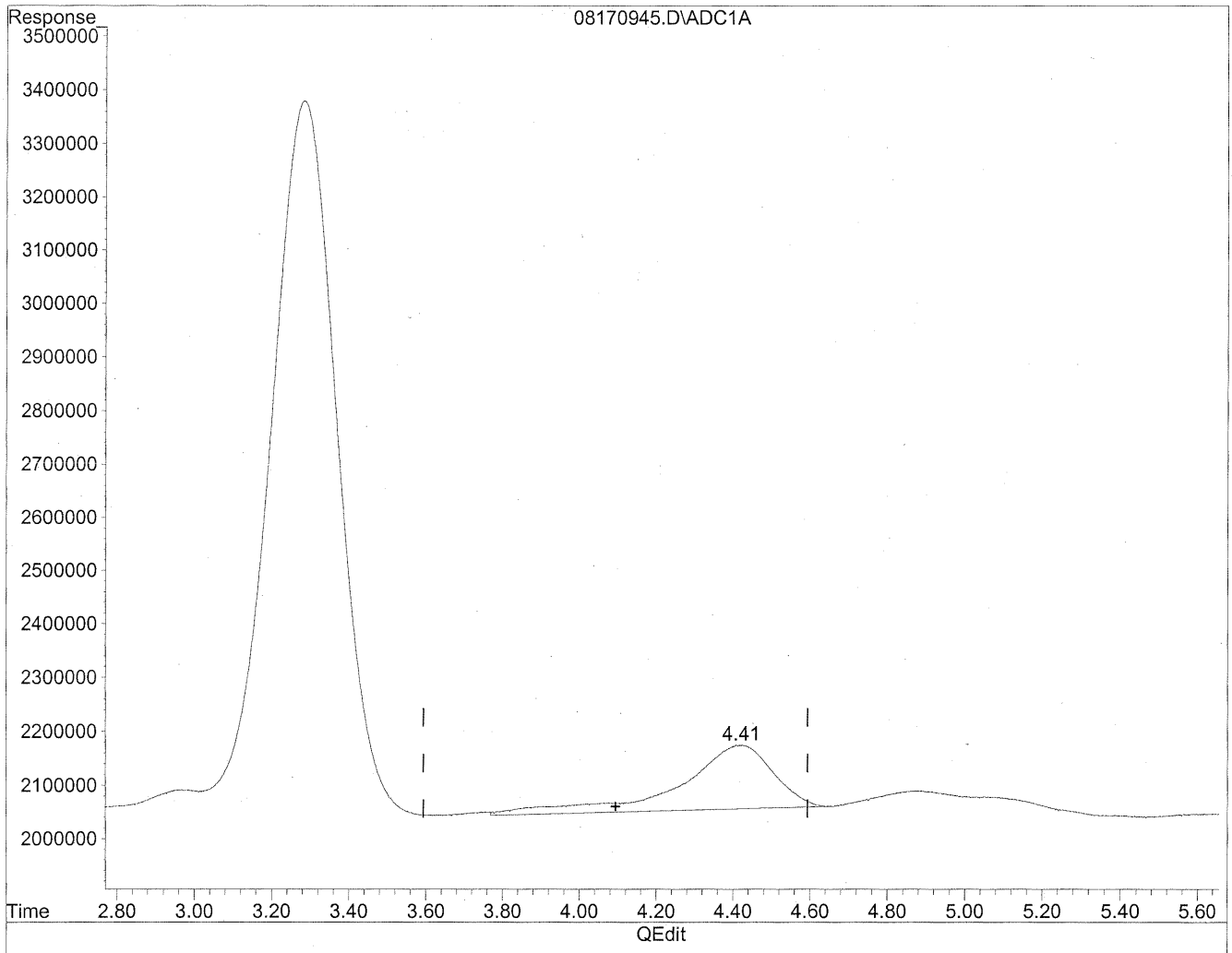
318

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170945.D Vial: 44
Acq On : 18 Aug 2009 1:52 am Operator: HC
Sample : P0902772-012 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:36 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.42min 207.774ng/ml

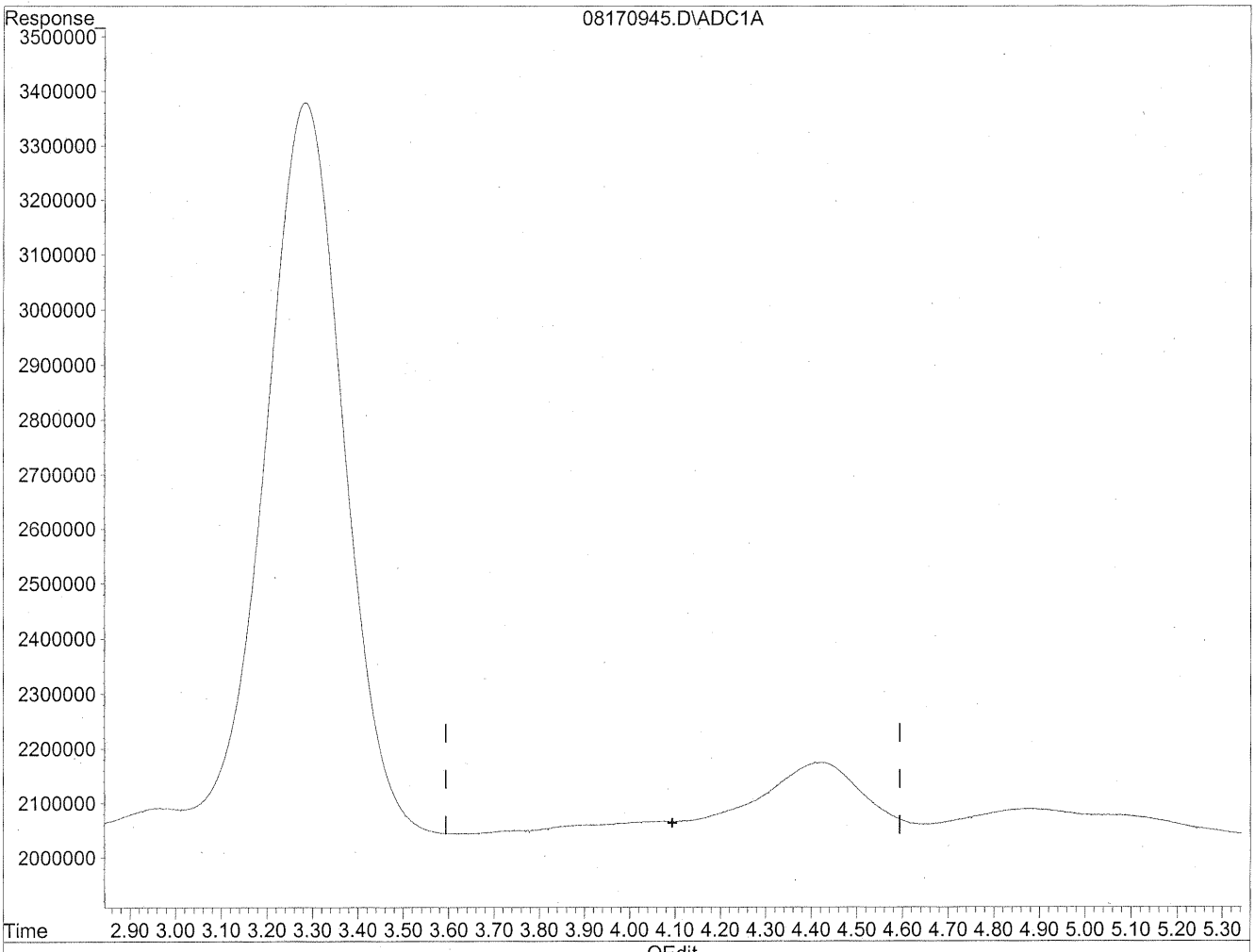
response 20240383

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170945.D Vial: 44
Acq On : 18 Aug 2009 1:52 am Operator: HC
Sample : P0902772-012 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 12:36 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

htc 8/22/09 mrp

KP 8/23/09

320

(+) = Expected Retention Time

08170945.D TO110709.M Sat Aug 22 12:36:58 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **Method Blank**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P090817-MB

Test Code: EPA Method TO-11A
Instrument ID: Waters LC Module I Plus/UV_Vis 360/LC1
Analyst: Hani Cherazaie
Sampling Media: Silica Gel DNPH Tube
Test Notes: **BC**

Date Collected: NA
Date Received: NA
Date Analyzed: 08/17/09
Desorption Volume: 1.0 ml
Volume Sampled: NA Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	Formaldehyde	< 100	NA	NA	NA	NA	
75-07-0	Acetaldehyde	< 100	NA	NA	NA	NA	
123-38-6	Propionaldehyde	< 100	NA	NA	NA	NA	
4170-30-3	Crotonaldehyde, Total	< 100	NA	NA	NA	NA	
123-72-8	Butyraldehyde	< 100	NA	NA	NA	NA	
100-52-7	Benzaldehyde	< 100	NA	NA	NA	NA	
590-86-3	Isovaleraldehyde	< 100	NA	NA	NA	NA	
110-62-3	Valeraldehyde	< 100	NA	NA	NA	NA	
529-20-4	o-Tolualdehyde	< 100	NA	NA	NA	NA	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	NA	NA	NA	NA	
66-25-1	n-Hexaldehyde	< 100	NA	NA	NA	NA	
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	NA	NA	NA	NA	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
NA = Not applicable.

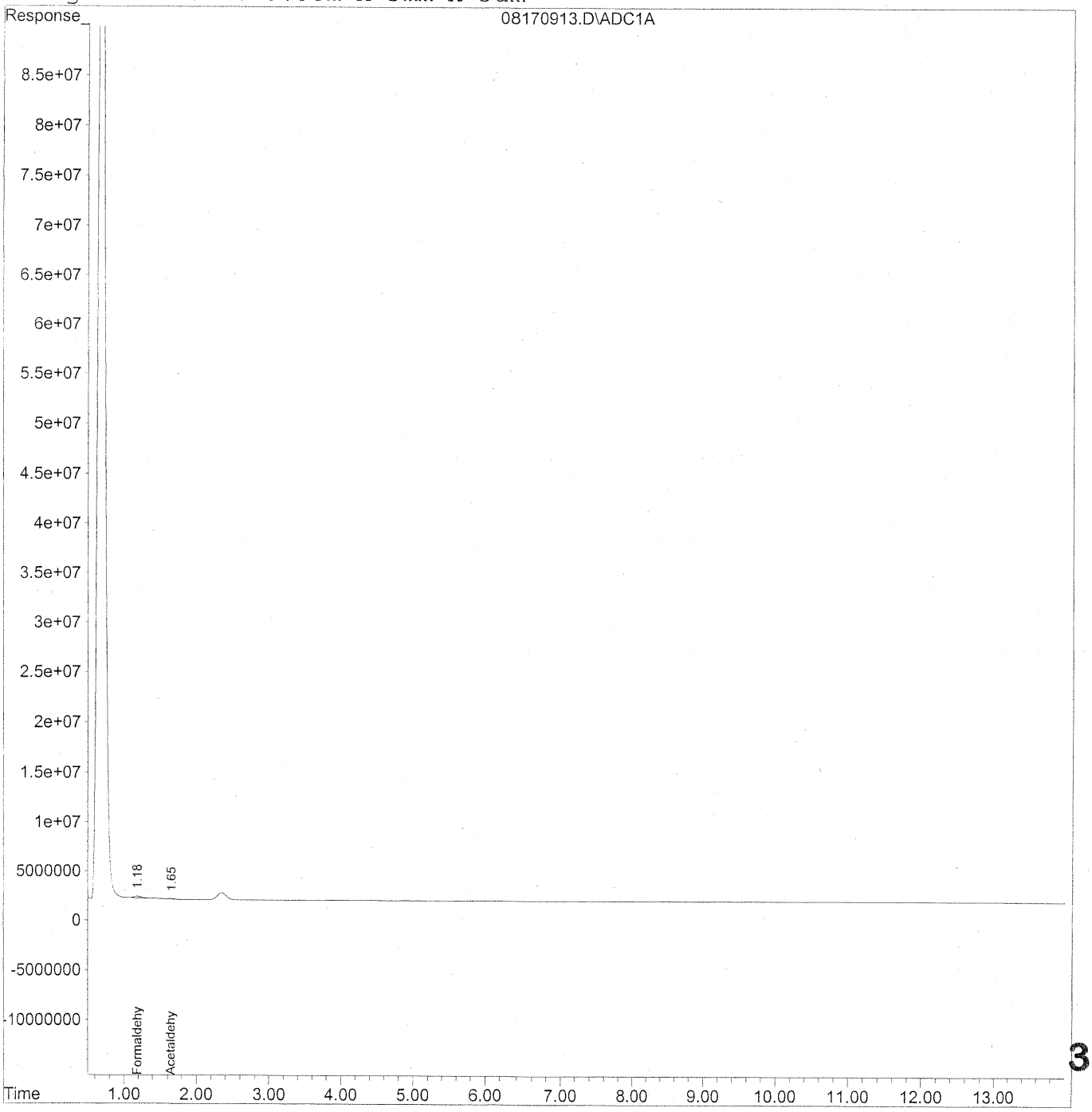
Verified By: _____ Date: 8/25/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170913.D Vial: 13
Acq On : 17 Aug 2009 5:50 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 20 17:30 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



322

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170913.D Vial: 13
Acq On : 17 Aug 2009 5:50 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 20 17:30 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

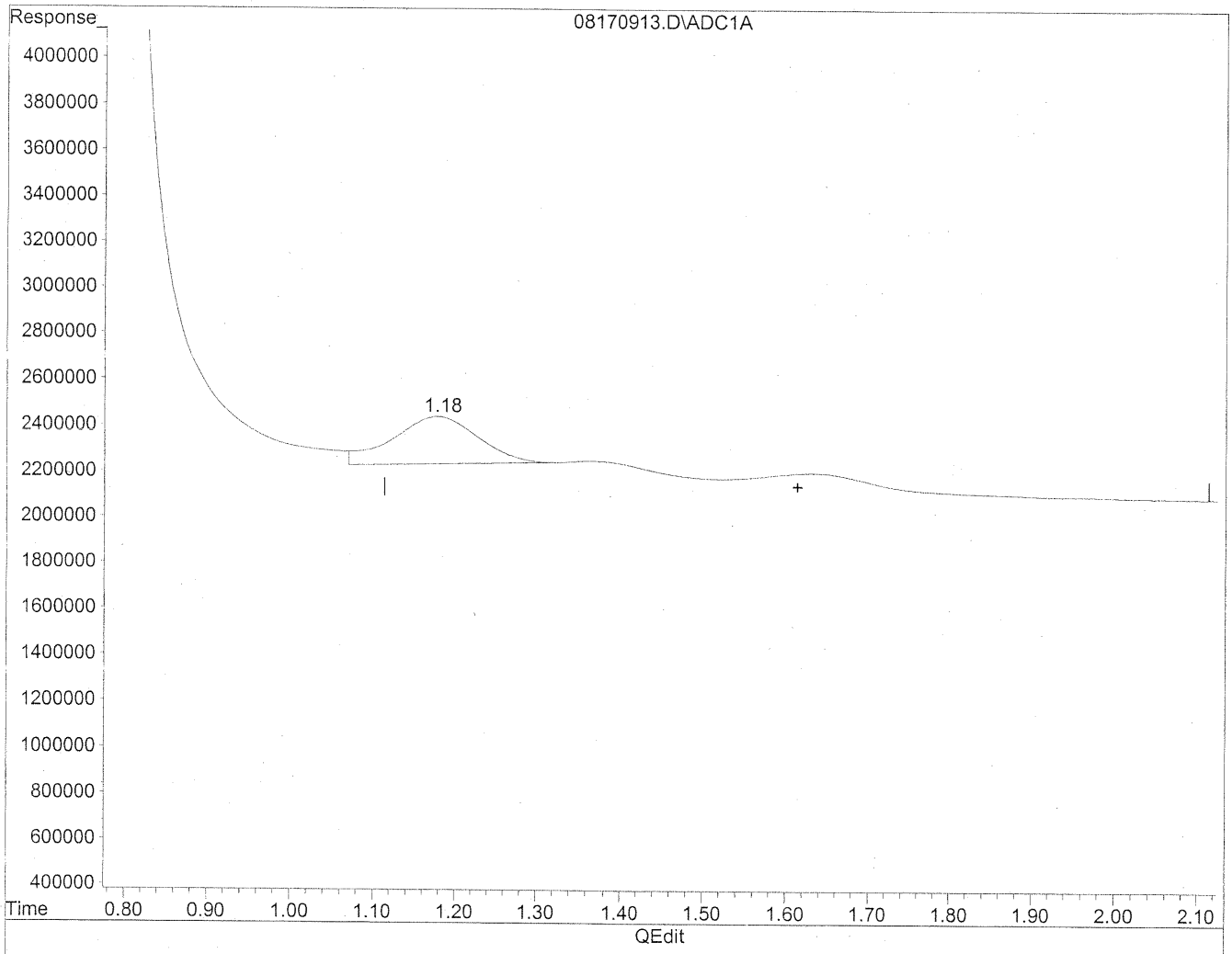
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.18	10839190	59.043	ng/mlm
2)	Acetaldehyde	1.65	4913326	35.039	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170913.D Vial: 13
Acq On : 17 Aug 2009 5:50 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 20 17:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.18min 106.863ng/ml

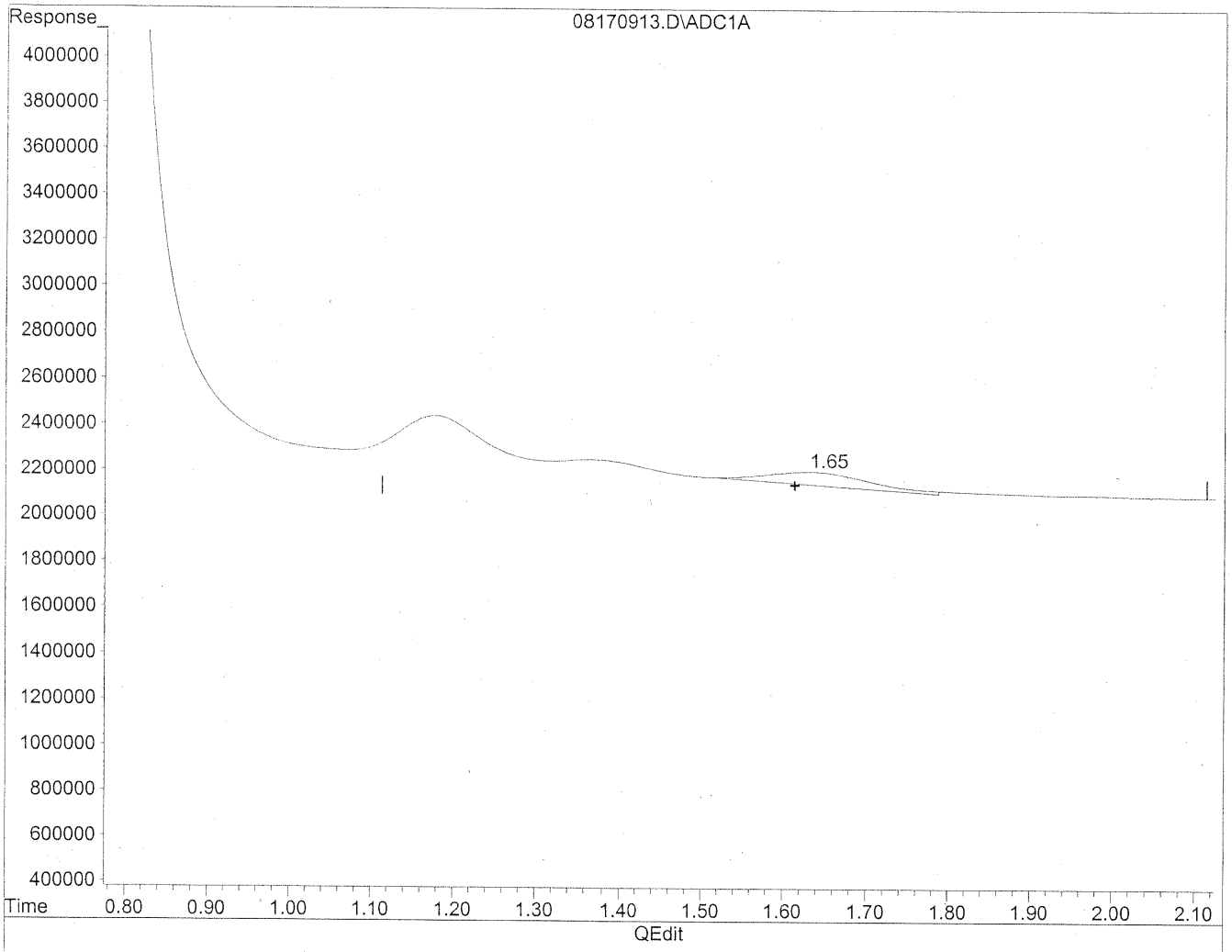
response 14984773

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170913.D Vial: 13
Acq On : 17 Aug 2009 5:50 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 20 17:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.65min 35.039ng/ml m

response 4913326

HC 8/21/09
WP 8/25/09

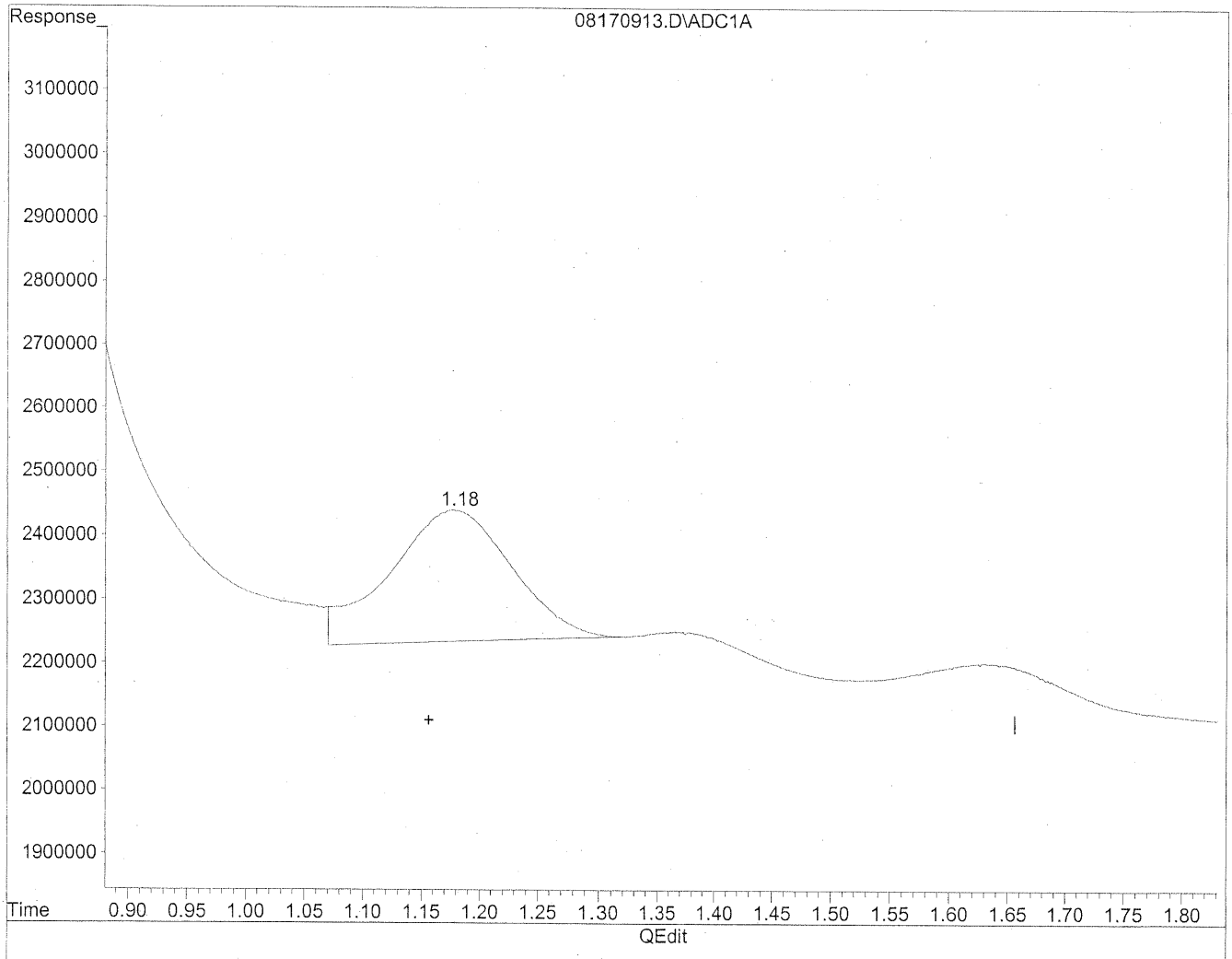
325

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170913.D Vial: 13
Acq On : 17 Aug 2009 5:50 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 20 17:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.18min 81.625ng/ml

response 14984773

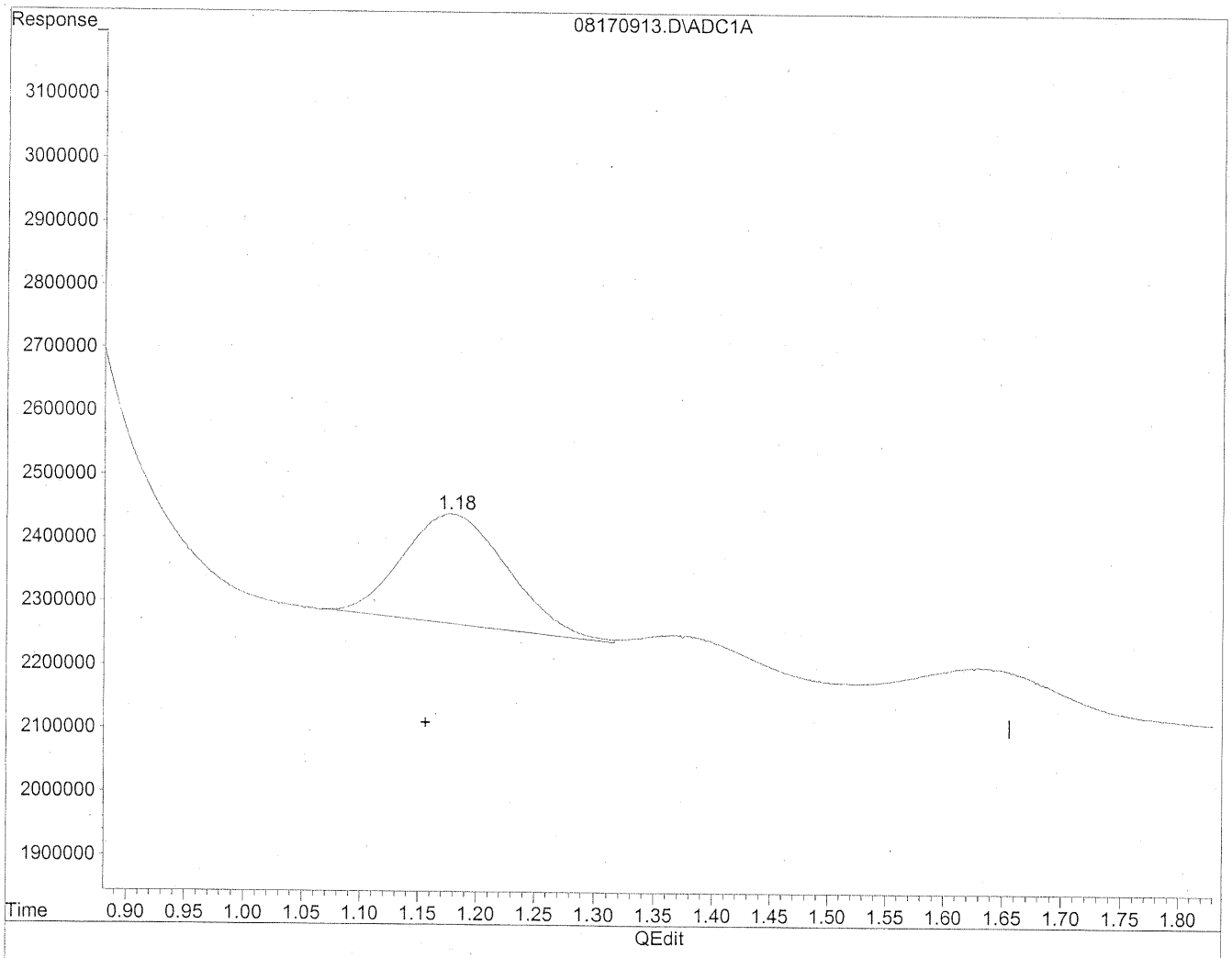
326

(+) = Expected Retention Time
08170913.D TO110709.M Thu Aug 20 17:30:30 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170913.D Vial: 13
Acq On : 17 Aug 2009 5:50 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 20 17:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.18min 59.043ng/ml m

response 10839190

HC 8/21/09
IC
KP 8/23/09

327

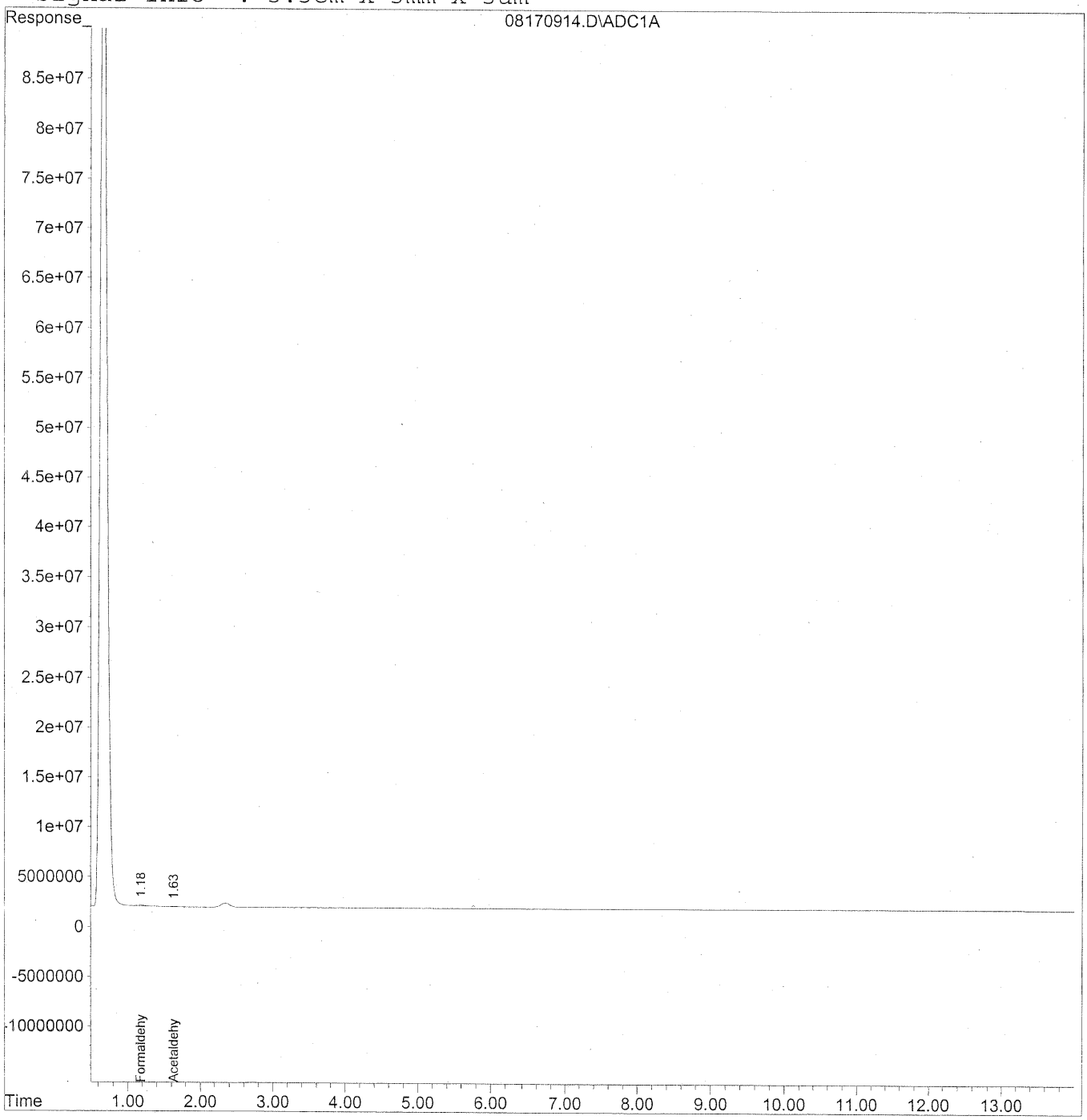
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170914.D Vial: 14
Acq On : 17 Aug 2009 6:05 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 20 17:36 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



328

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170914.D Vial: 14
Acq On : 17 Aug 2009 6:05 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 20 17:36 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

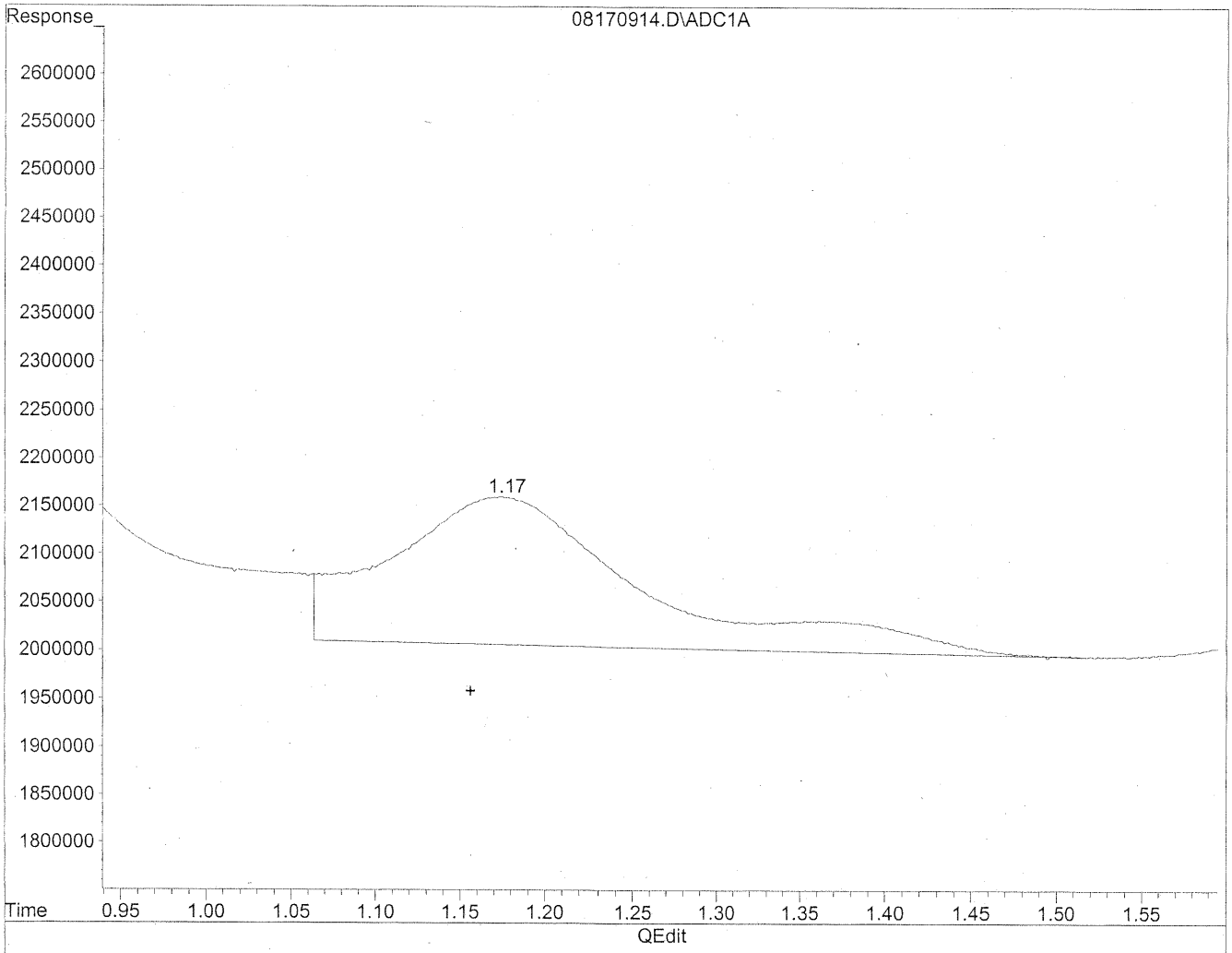
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.18	6589179	35.892	ng/mlm
2)	Acetaldehyde	1.63	2141330	15.271	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170914.D Vial: 14
Acq On : 17 Aug 2009 6:05 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 20 17:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 86.671ng/ml

response 15911145

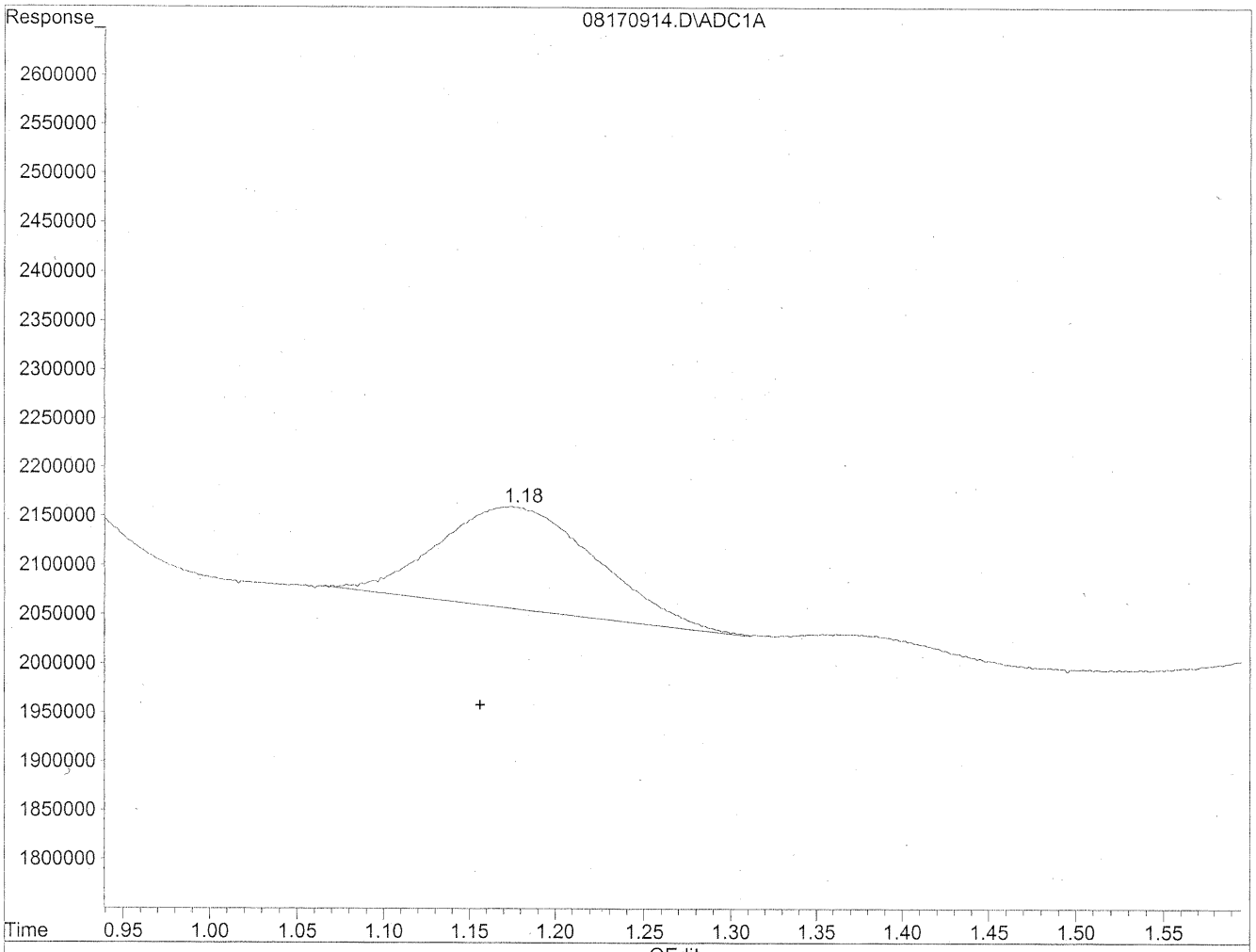
330

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170914.D Vial: 14
Acq On : 17 Aug 2009 6:05 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 20 17:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.18min 35.892ng/ml m

response 6589179

HC 8/24/09

IC 8/23/09

331

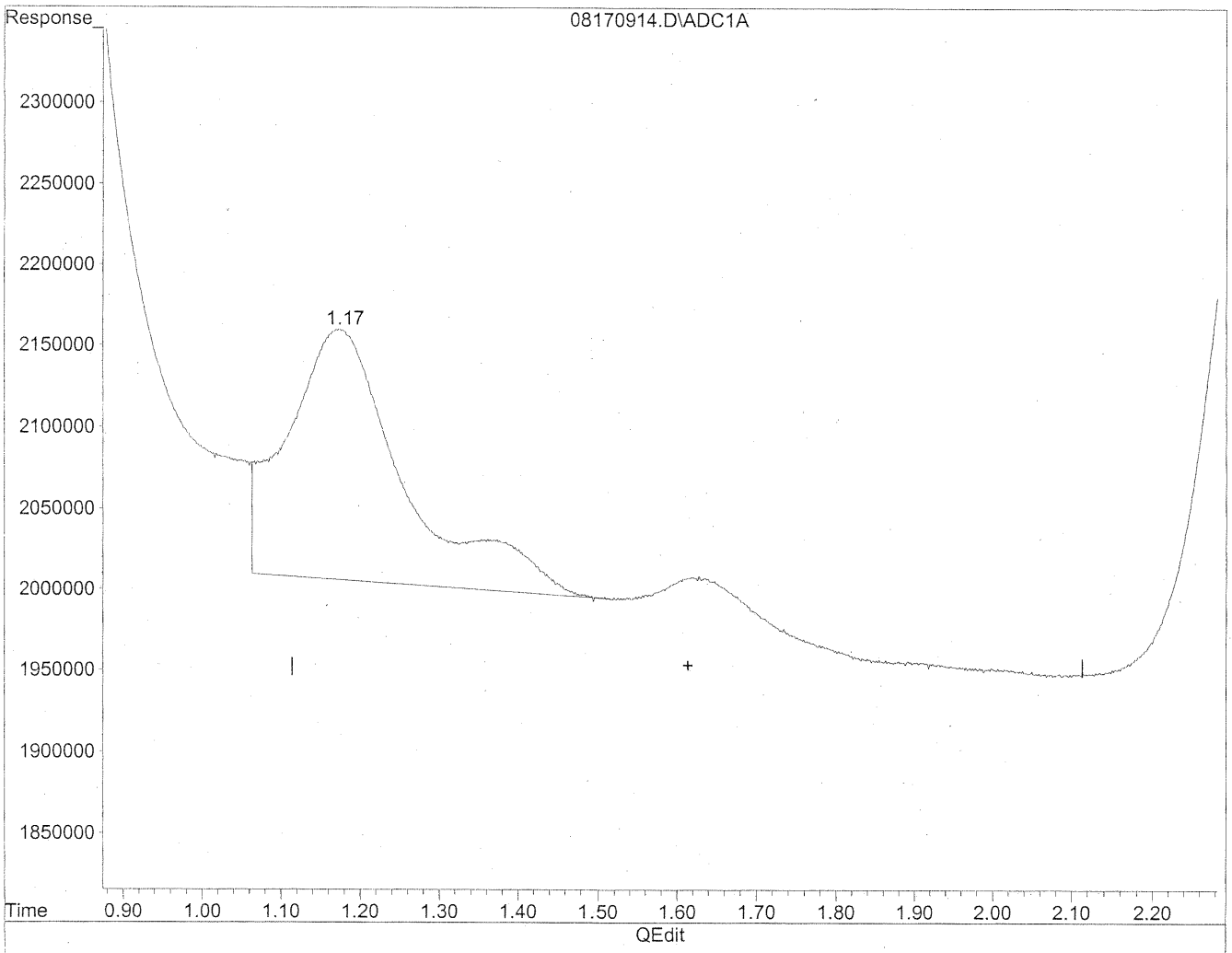
(+) = Expected Retention Time

08170914.D TO110709.M Thu Aug 20 17:36:11 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170914.D Vial: 14
Acq On : 17 Aug 2009 6:05 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 20 17:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.17min 113.470ng/ml

response 15911145

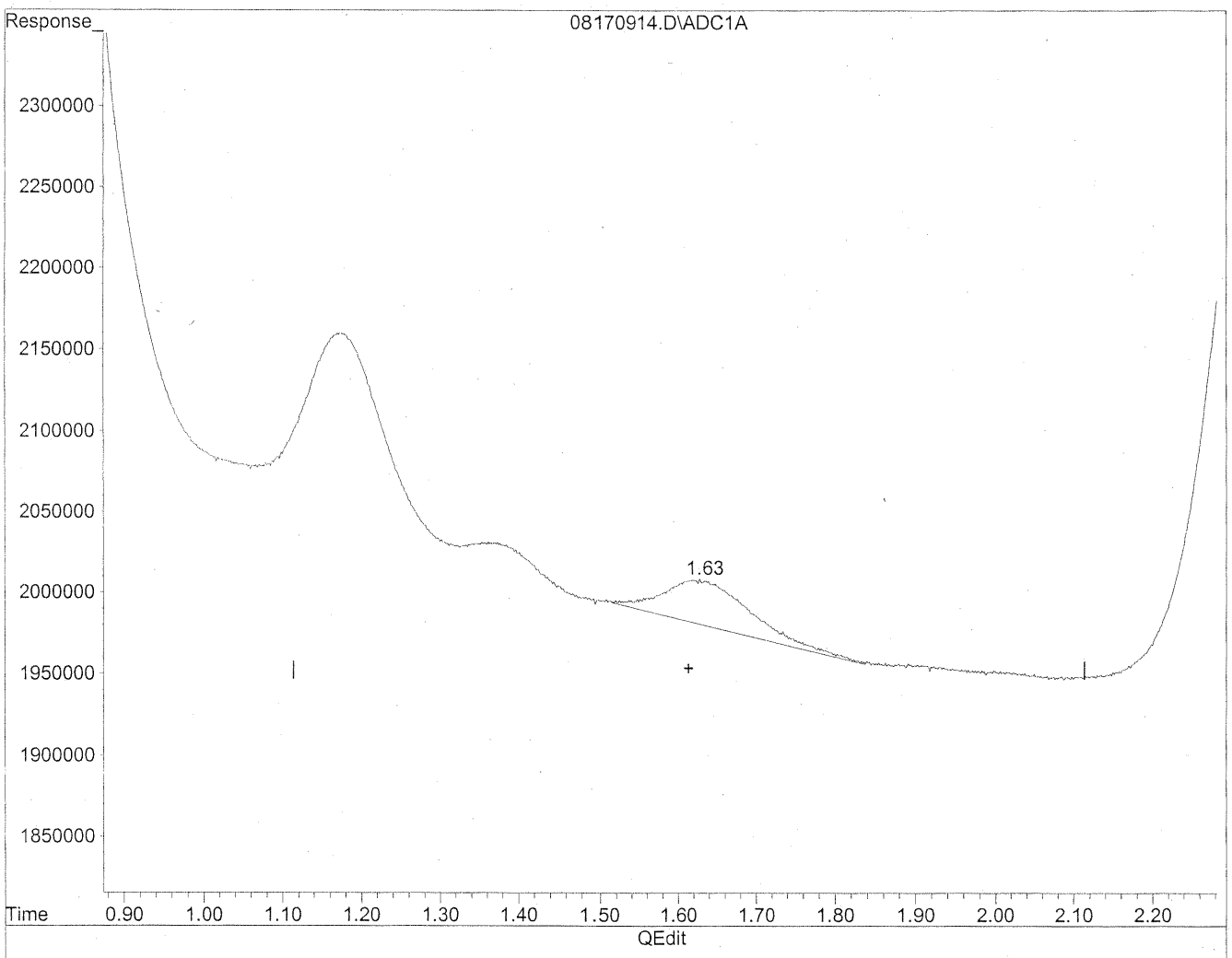
332

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170914.D Vial: 14
Acq On : 17 Aug 2009 6:05 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 20 17:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.63min 15.271ng/ml m

response 2141330

HC 8/21/09

1C

KB 8/23/09

333

(+) = Expected Retention Time

08170914.D TO110709.M Thu Aug 20 17:36:34 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **Method Blank (00:06)**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P090818-MB

Test Code: EPA Method TO-11A
Instrument ID: Waters LC Module I Plus/UV_Vis 360/LC1
Analyst: Hani Cherazaie
Sampling Media: Silica Gel DNPH Tube
Test Notes: **BC**

Date Collected: NA
Date Received: NA
Date Analyzed: 08/18/09
Desorption Volume: 1.0 ml
Volume Sampled: NA Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	Formaldehyde	< 100	NA	NA	NA	NA	
75-07-0	Acetaldehyde	< 100	NA	NA	NA	NA	
123-38-6	Propionaldehyde	< 100	NA	NA	NA	NA	
4170-30-3	Crotonaldehyde, Total	< 100	NA	NA	NA	NA	
123-72-8	Butyraldehyde	< 100	NA	NA	NA	NA	
100-52-7	Benzaldehyde	< 100	NA	NA	NA	NA	
590-86-3	Isovaleraldehyde	< 100	NA	NA	NA	NA	
110-62-3	Valeraldehyde	< 100	NA	NA	NA	NA	
529-20-4	o-Tolualdehyde	< 100	NA	NA	NA	NA	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	NA	NA	NA	NA	
66-25-1	n-Hexaldehyde	< 100	NA	NA	NA	NA	
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	NA	NA	NA	NA	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
NA = Not applicable.

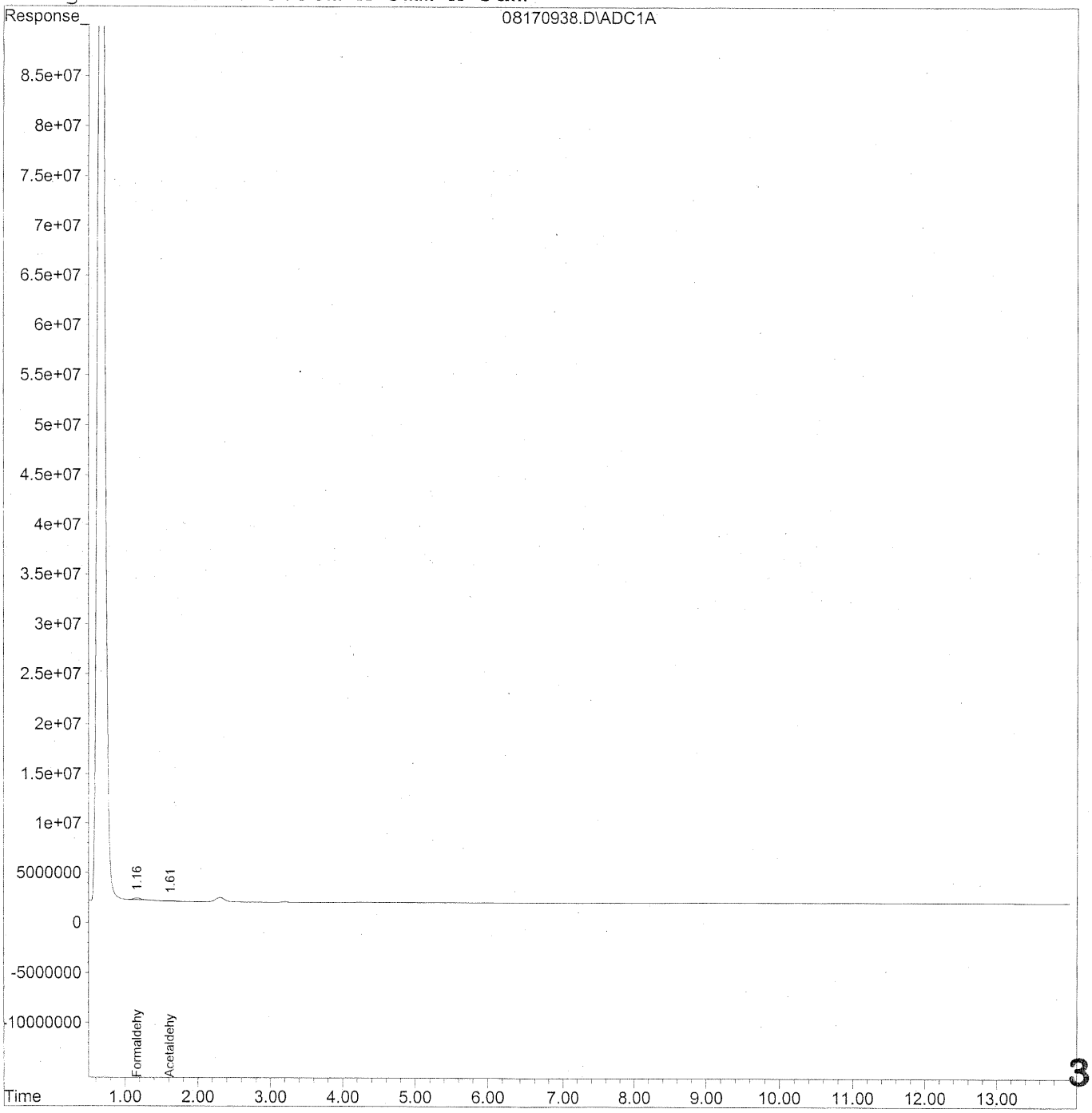
Verified By: _____ Date: 8/25/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170938.D Vial: 37
Acq On : 18 Aug 2009 12:06 am Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:17 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



335

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170938.D Vial: 37
Acq On : 18 Aug 2009 12:06 am Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:17 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

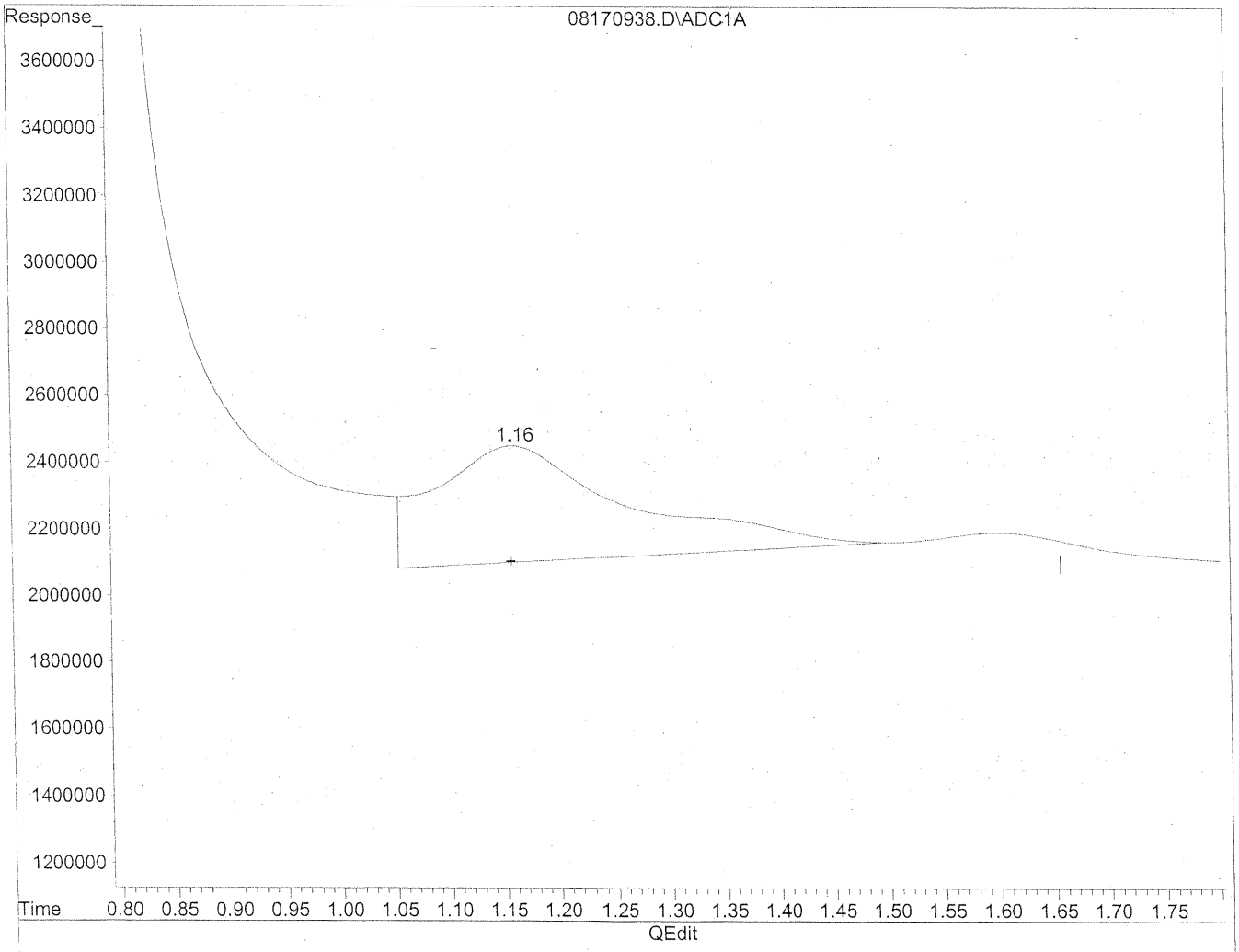
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	11165898	60.823	ng/mlm
2)	Acetaldehyde	1.61	3134248	22.352	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170938.D Vial: 37
Acq On : 18 Aug 2009 12:06 am Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:17 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.16min 233.224ng/ml

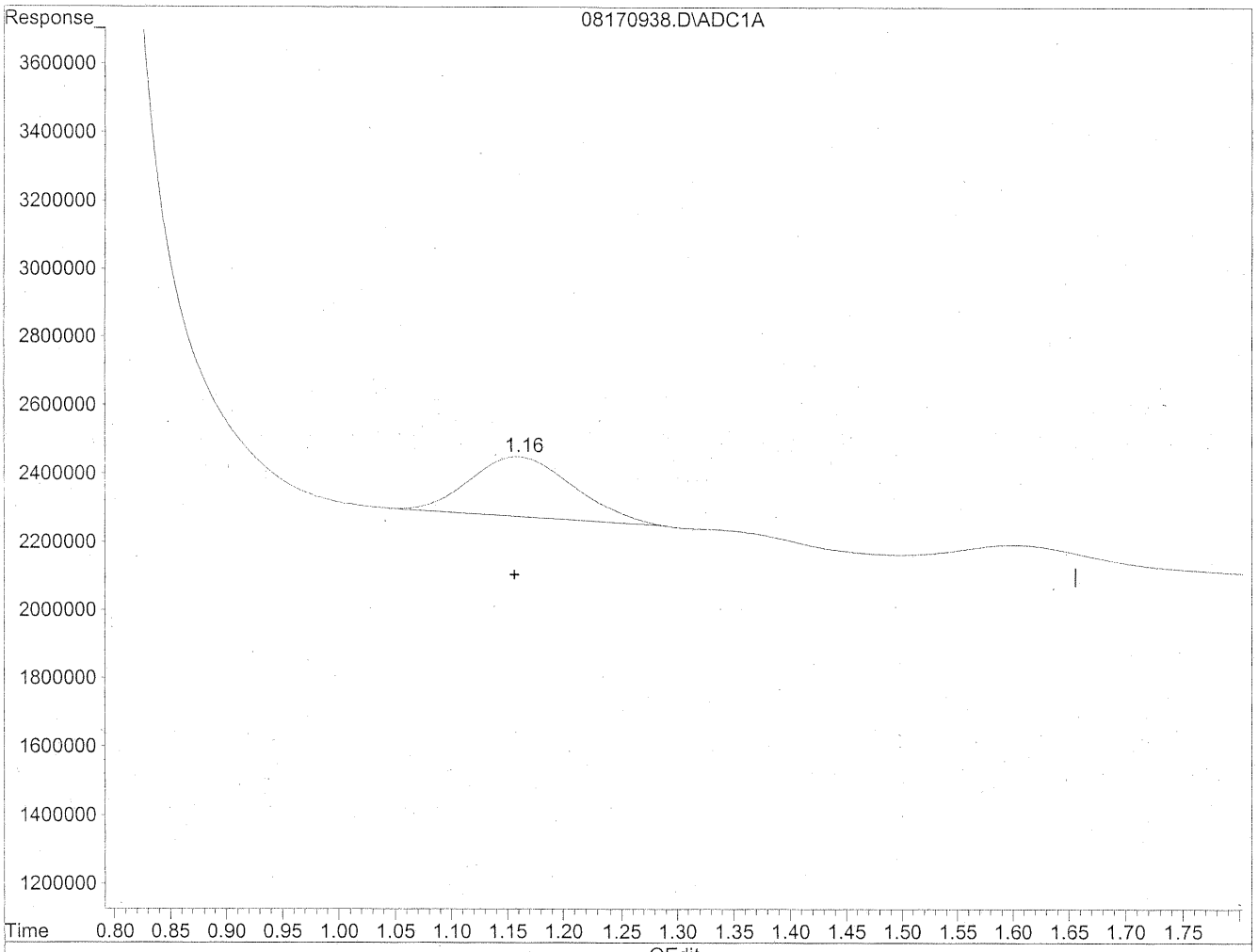
response 42815580

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170938.D Vial: 37
Acq On : 18 Aug 2009 12:06 am Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:17 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.16min 60.823ng/ml m

response 11165898

HC 8/21/09 LC

KPB 8/23/09

338

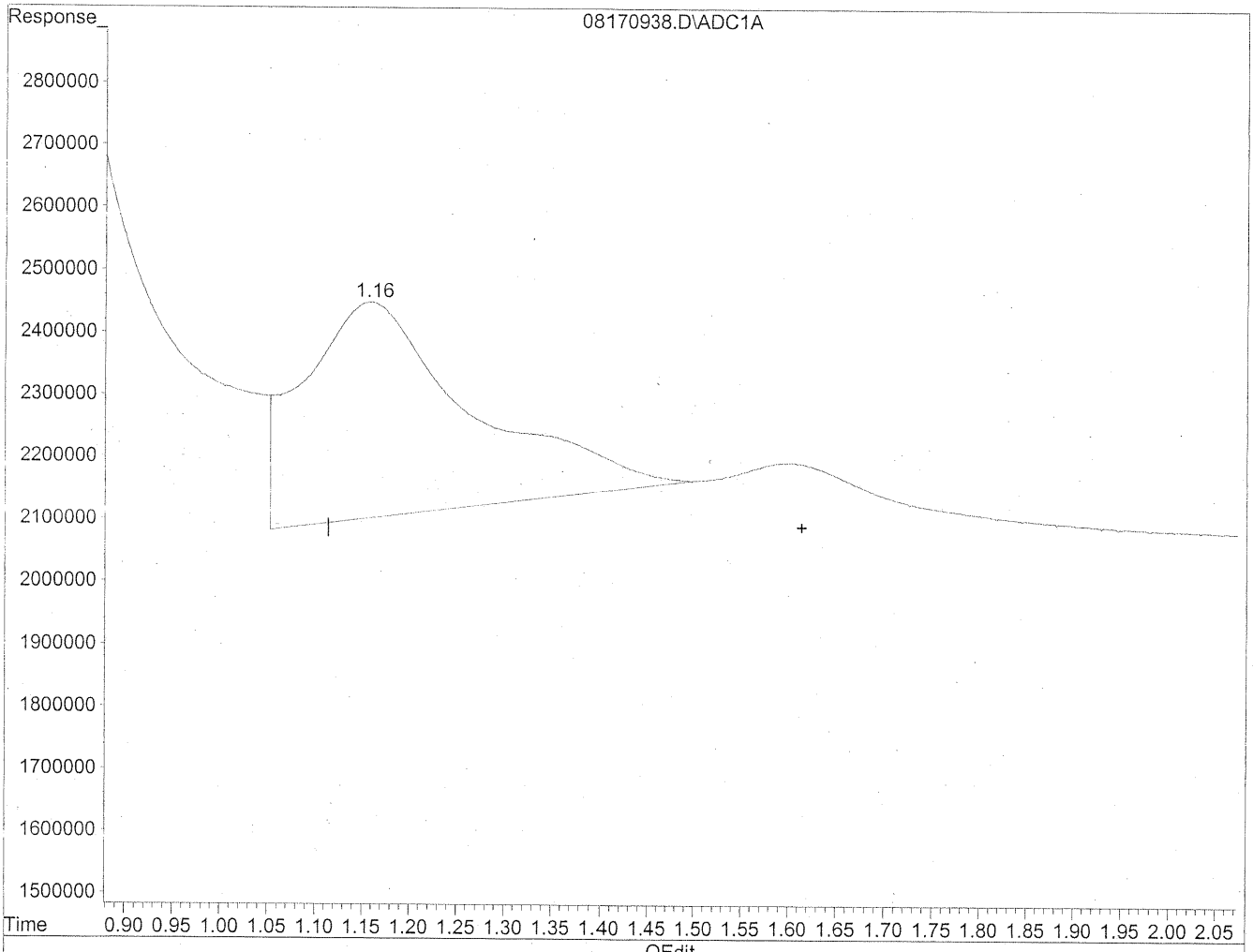
(+) = Expected Retention Time

08170938.D TO110709.M Fri Aug 21 18:17:27 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170938.D Vial: 37
Acq On : 18 Aug 2009 12:06 am Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:17 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(2) Acetaldehyde

1.16min 305.338ng/ml

response 42815580

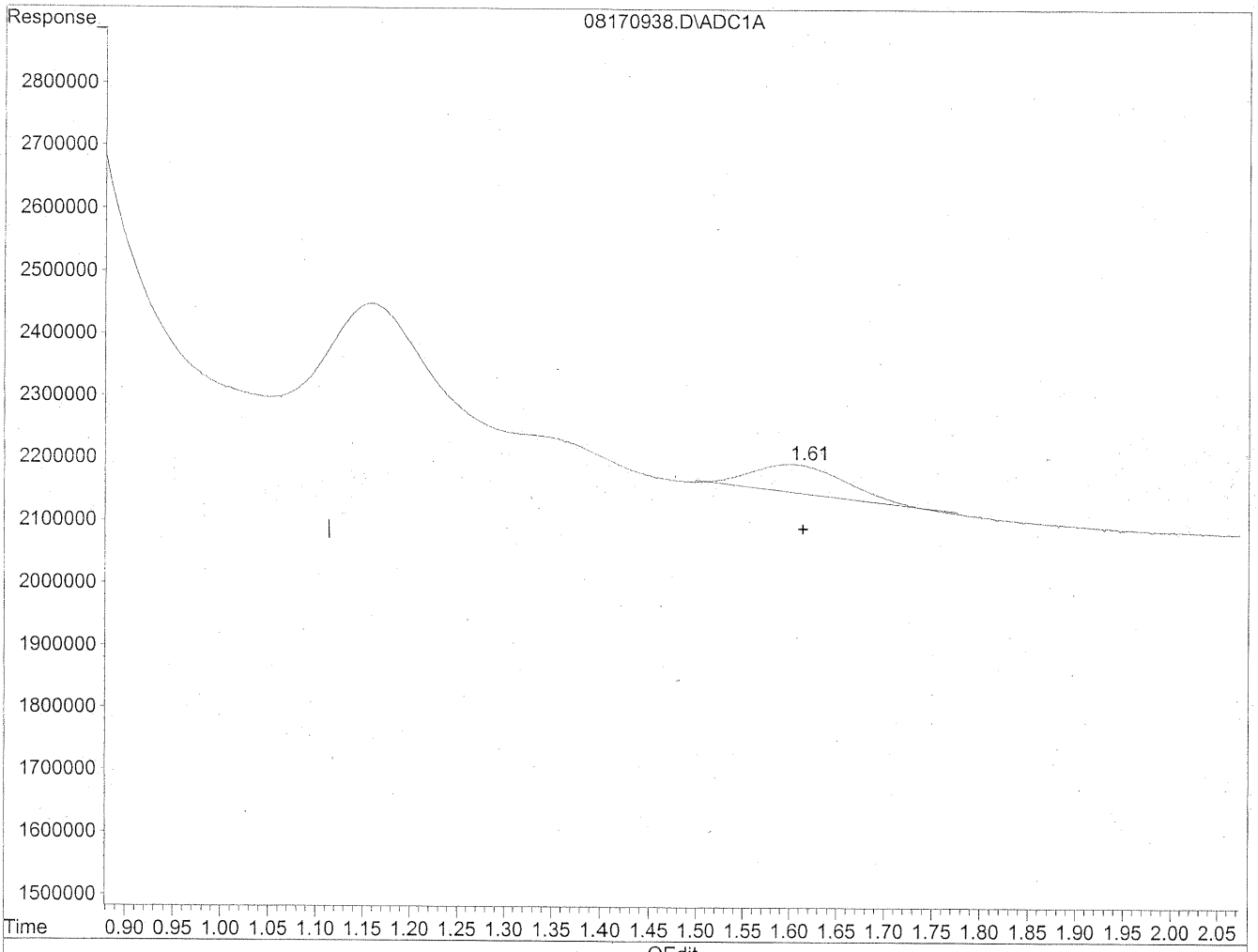
(+) = Expected Retention Time

08170938.D TO110709.M Fri Aug 21 18:17:35 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170938.D Vial: 37
Acq On : 18 Aug 2009 12:06 am Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:17 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.61min 22.352ng/ml m

response 3134248

HC 8/21/09 LC

KP 8/23/09

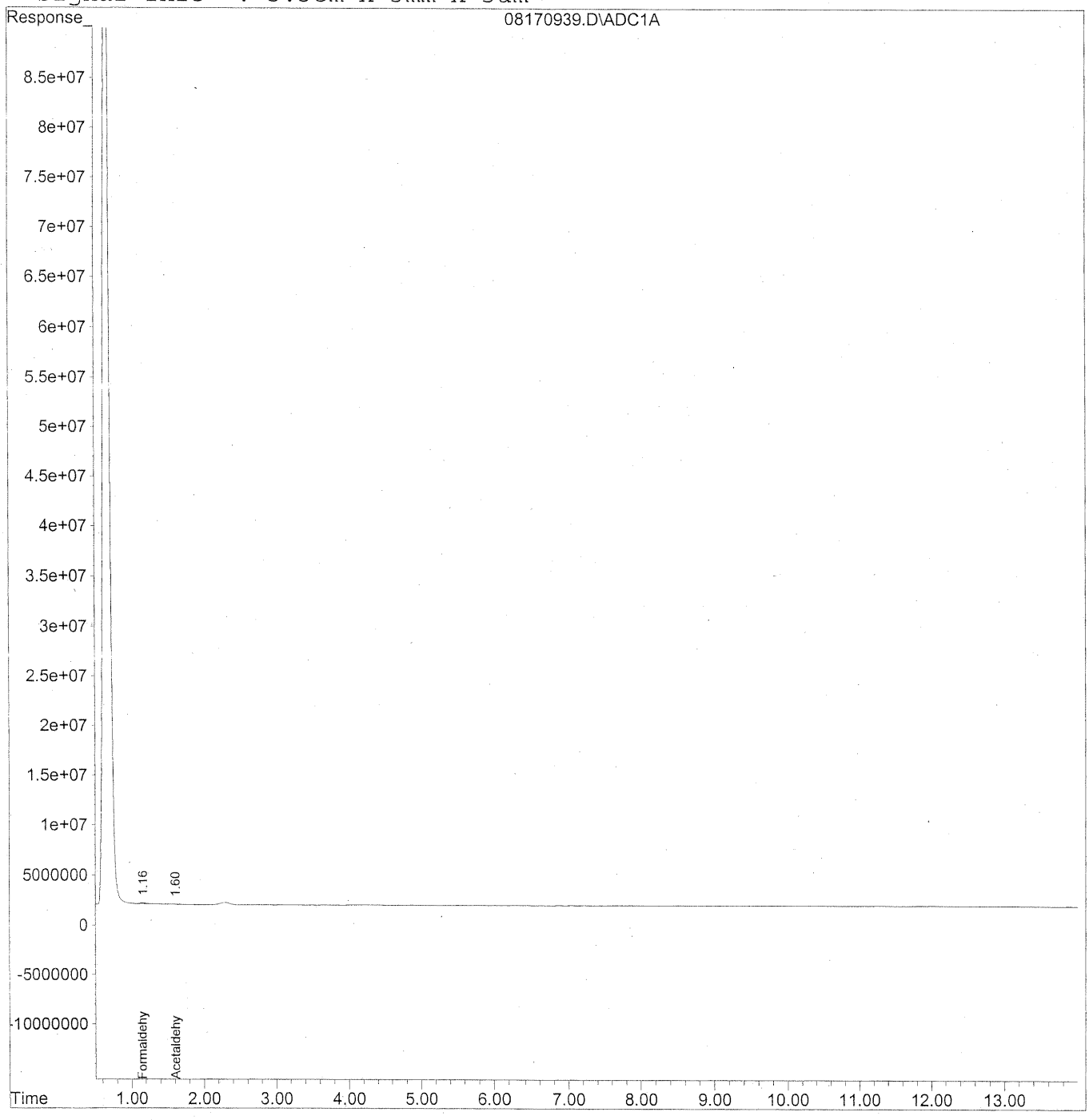
340

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_08\17\08170939.D Vial: 38
Acq On : 18 Aug 2009 12:21 am Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:18 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170939.D Vial: 38
Acq On : 18 Aug 2009 12:21 am Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:18 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

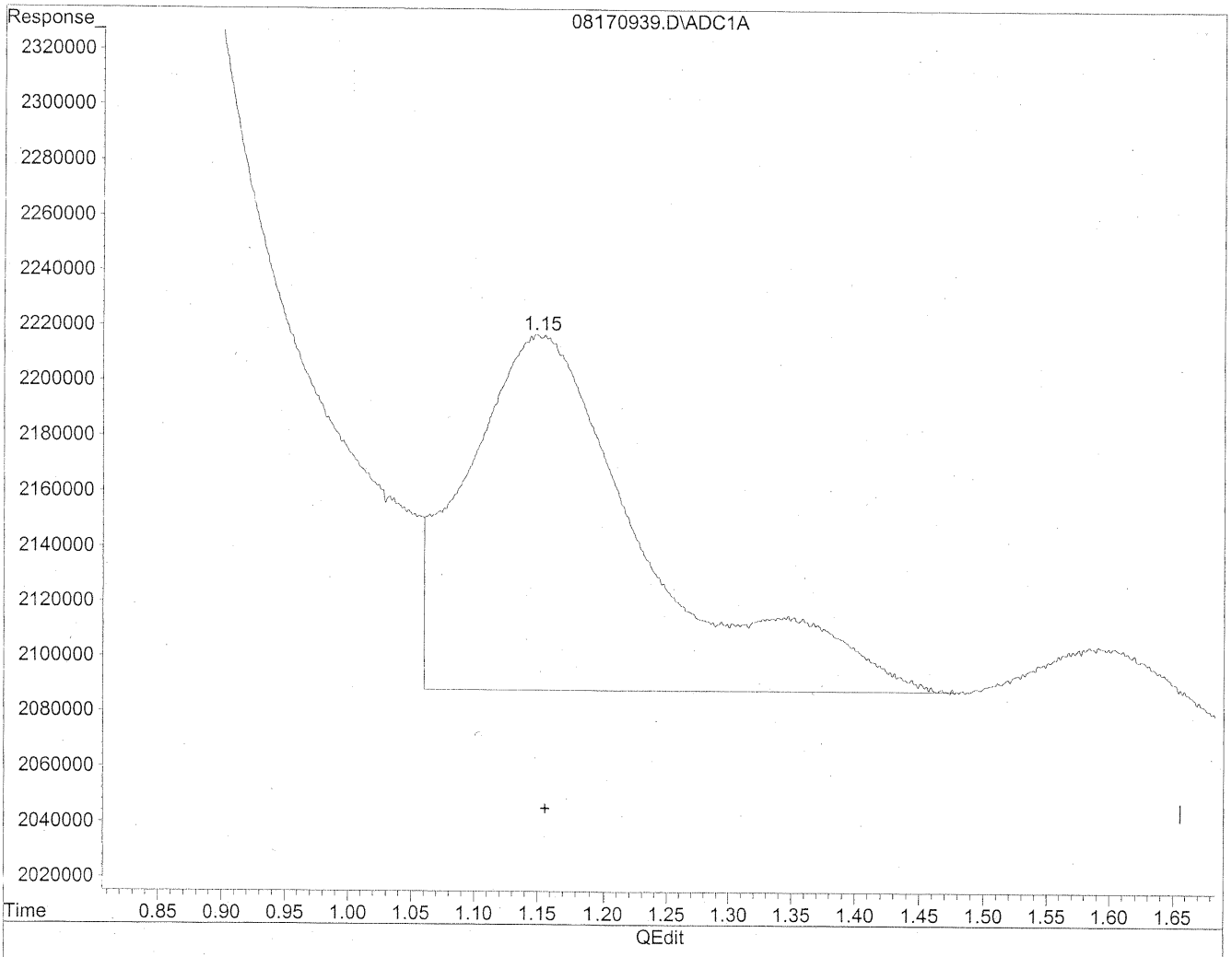
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	4863259	26.491	ng/mlm
2)	Acetaldehyde	1.60	2122505	15.137	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170939.D Vial: 38
Acq On : 18 Aug 2009 12:21 am Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:17 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.15min 69.012ng/ml

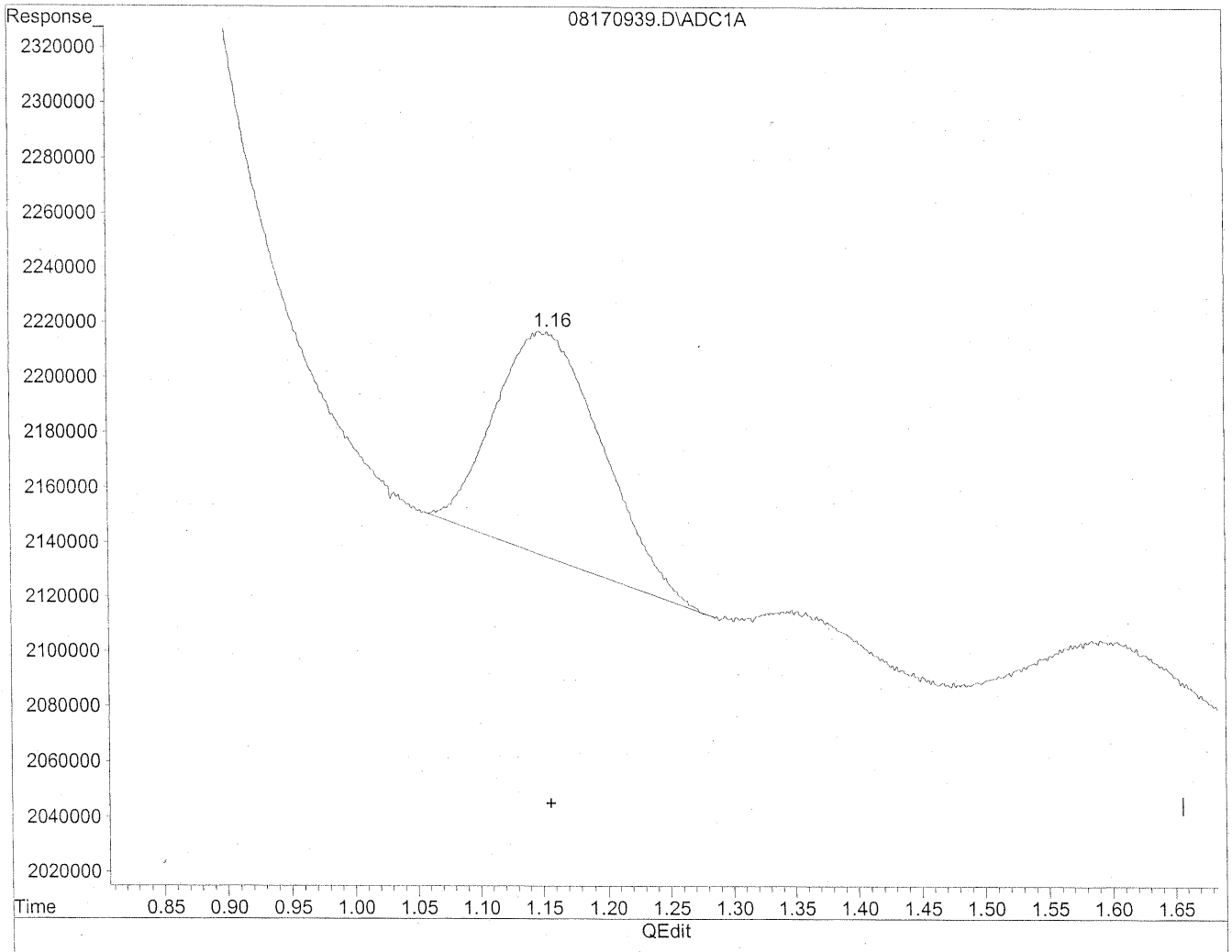
response 12669359

343

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_08\17\08170939.D Vial: 38
Acq On : 18 Aug 2009 12:21 am Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:17 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.16min 26.491ng/ml m

response 4863259

HC 8/21/09 LC

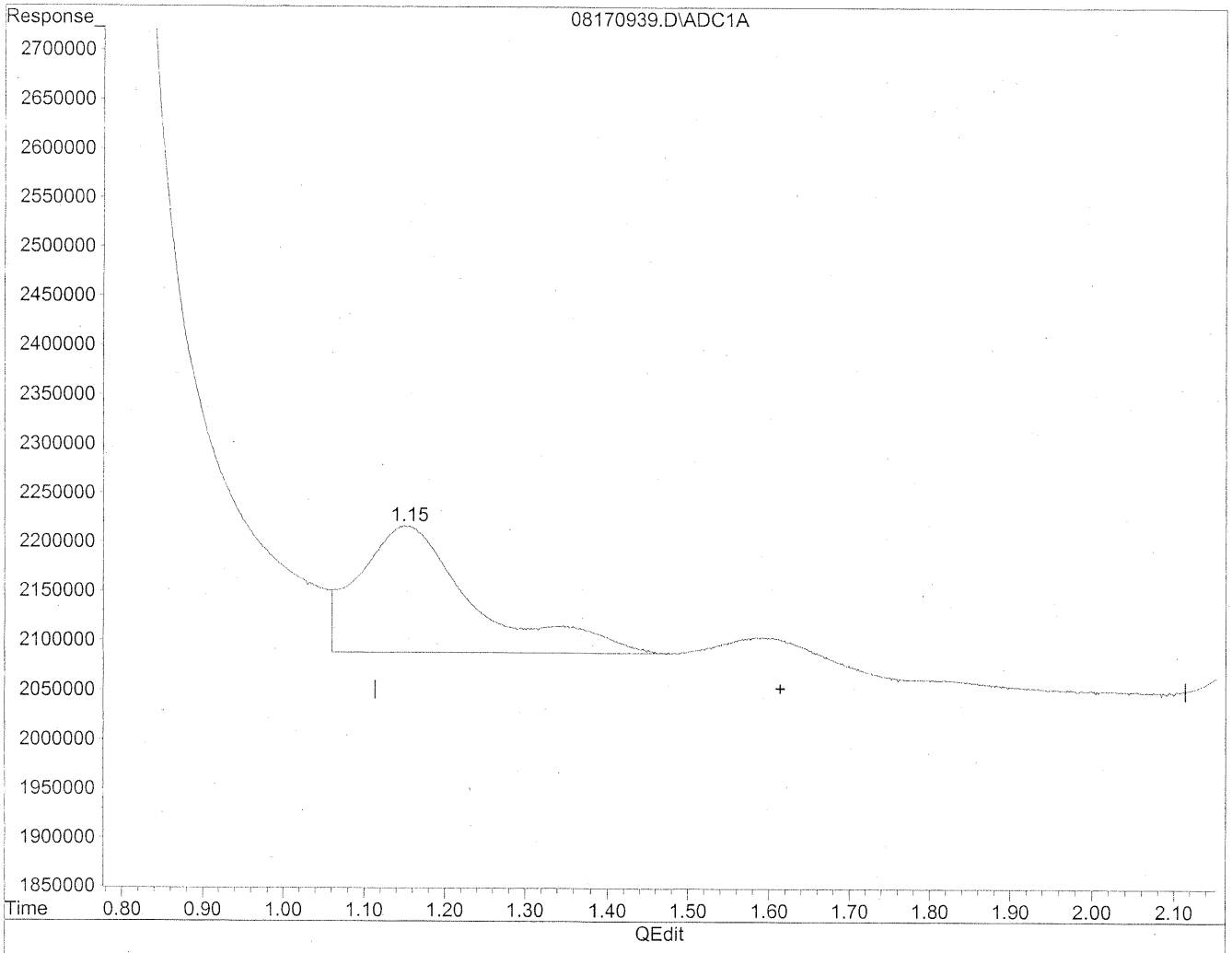
KPB 8/23/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170939.D Vial: 38
Acq On : 18 Aug 2009 12:21 am Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:17 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.15min 90.351ng/ml

response 12669359

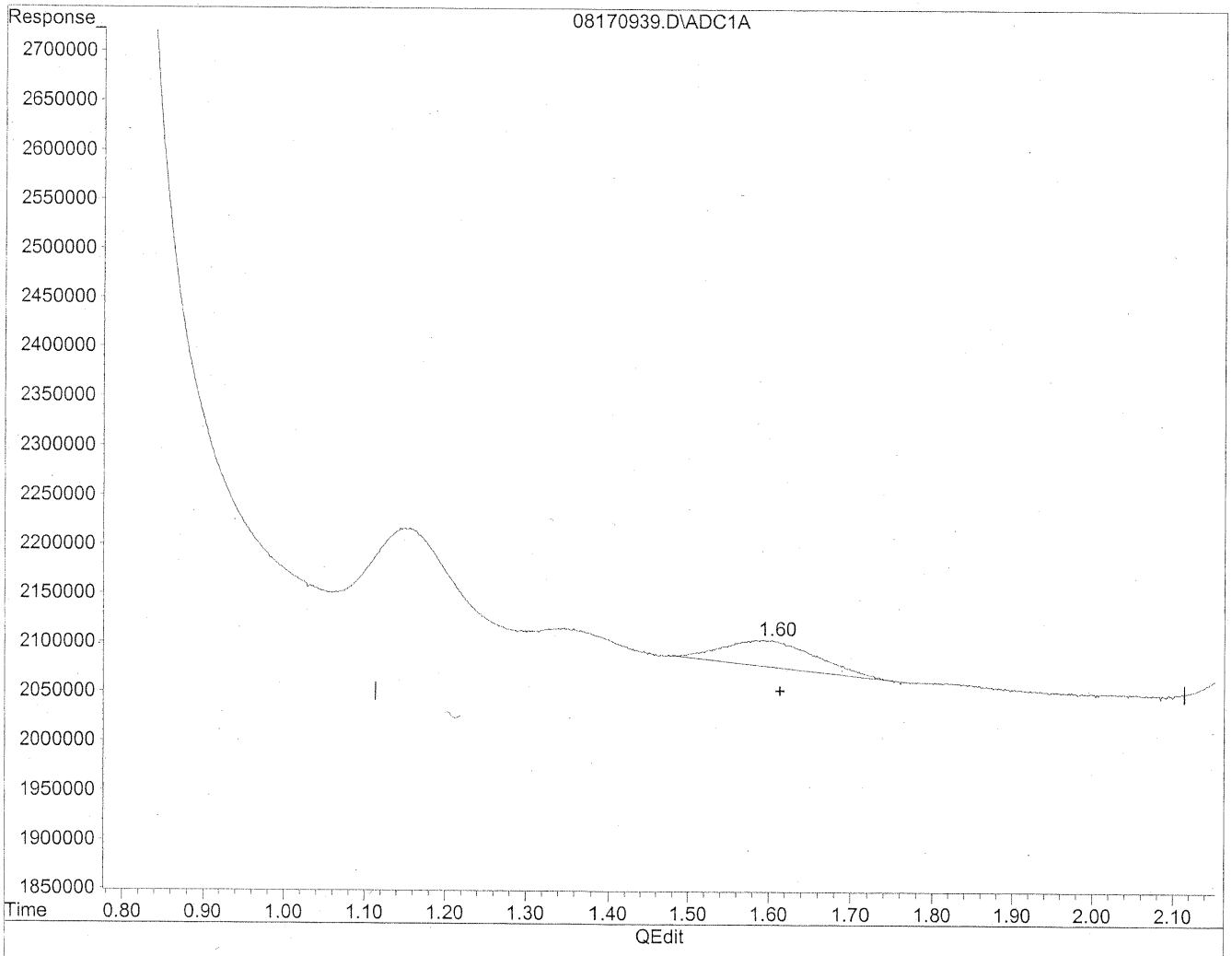
345

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170939.D Vial: 38
Acq On : 18 Aug 2009 12:21 am Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 21 18:17 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.60min 15.137ng/ml m

response 2122505

tlc 8/21/09 wp

8/23/09

(+) = Expected Retention Time

08170939.D TO110709.M Fri Aug 21 18:18:28 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **Method Blank (12:38)**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P090818-MB

Test Code: EPA Method TO-11A
Instrument ID: Waters LC Module I Plus/UV_Vis 360/LC1
Analyst: Hani Cherazaie
Sampling Media: Silica Gel DNPH Tube
Test Notes: **BC**

Date Collected: NA
Date Received: NA
Date Analyzed: 08/18/09
Desorption Volume: 1.0 ml
Volume Sampled: NA Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	Formaldehyde	< 100	NA	NA	NA	NA	
75-07-0	Acetaldehyde	< 100	NA	NA	NA	NA	
123-38-6	Propionaldehyde	< 100	NA	NA	NA	NA	
4170-30-3	Crotonaldehyde, Total	< 100	NA	NA	NA	NA	
123-72-8	Butyraldehyde	< 100	NA	NA	NA	NA	
100-52-7	Benzaldehyde	< 100	NA	NA	NA	NA	
590-86-3	Isovaleraldehyde	< 100	NA	NA	NA	NA	
110-62-3	Valeraldehyde	< 100	NA	NA	NA	NA	
529-20-4	o-Tolualdehyde	< 100	NA	NA	NA	NA	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	NA	NA	NA	NA	
66-25-1	n-Hexaldehyde	< 100	NA	NA	NA	NA	
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	NA	NA	NA	NA	

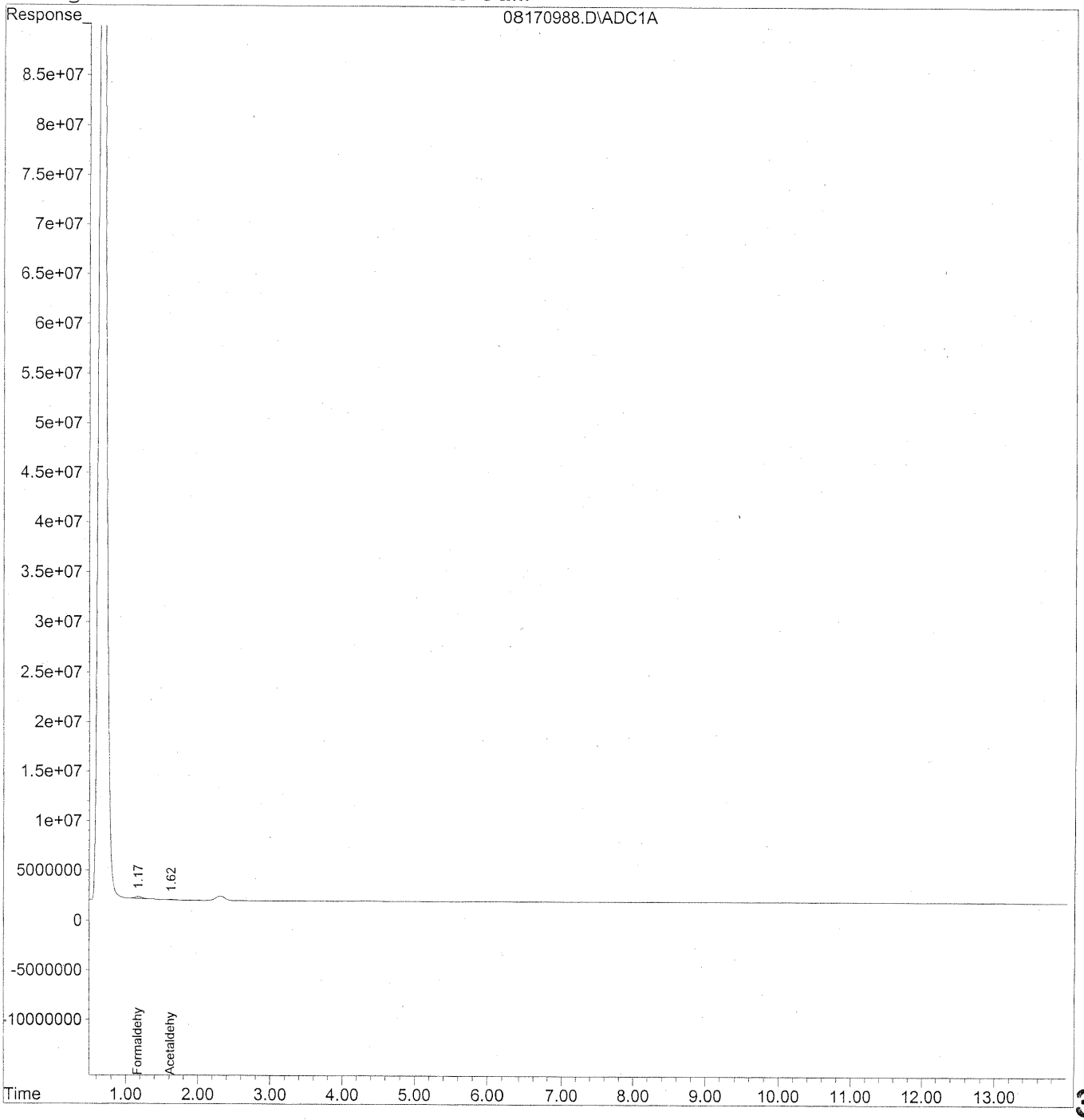
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
NA = Not applicable.

Verified By: _____ Date: 8/25/09

Data File : J:\LC01\DATA\TO11\2009_08\17\08170988.D Vial: 2
Acq On : 18 Aug 2009 12:38 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 13:17 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170988.D Vial: 2
Acq On : 18 Aug 2009 12:38 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 13:17 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

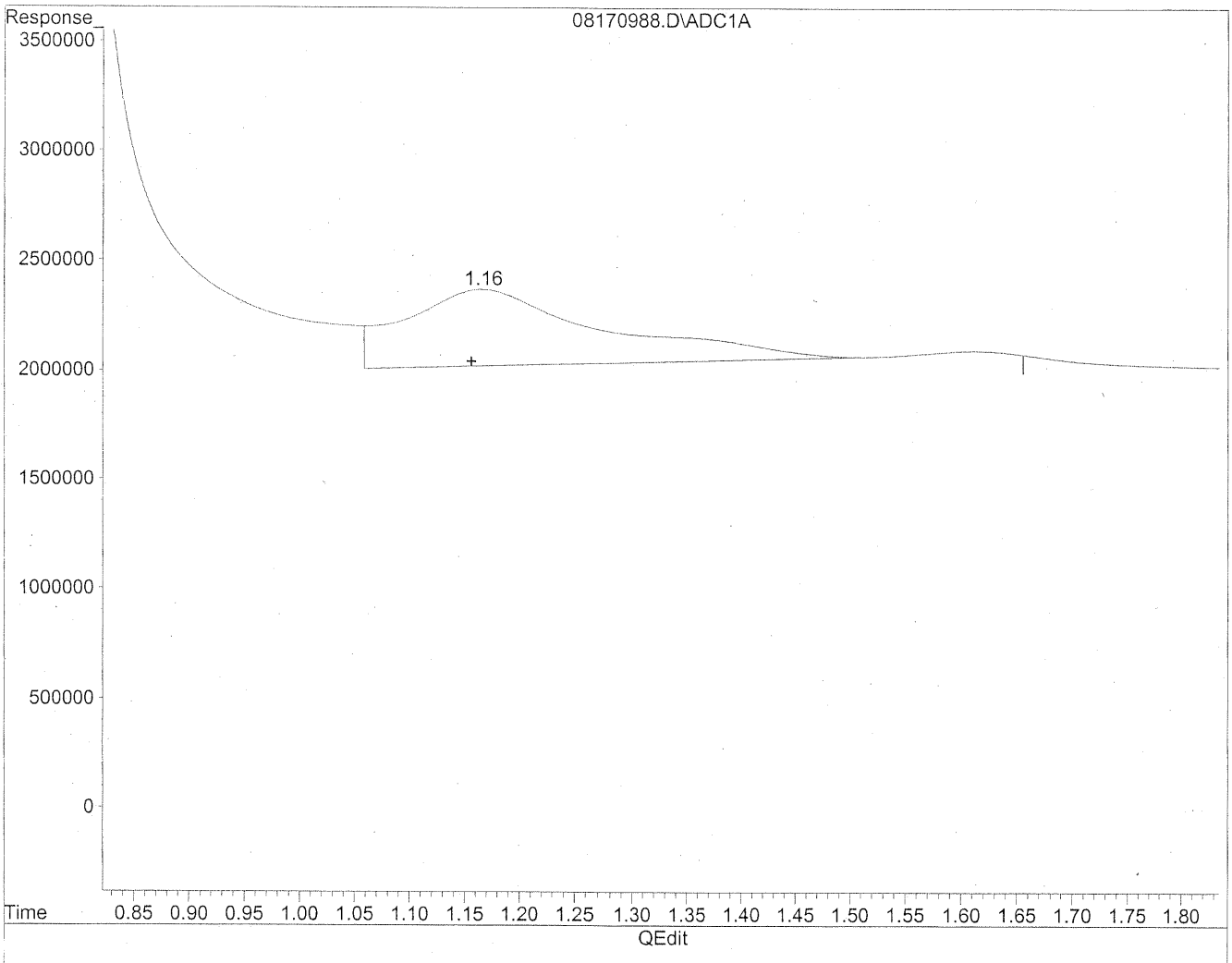
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
Target Compounds					
1)	Formaldehyde	1.17	12598459	68.626	ng/mlm
2)	Acetaldehyde	1.62	3332020	23.762	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170988.D Vial: 2
Acq On : 18 Aug 2009 12:38 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 13:16 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 235.655ng/ml

response 43261965

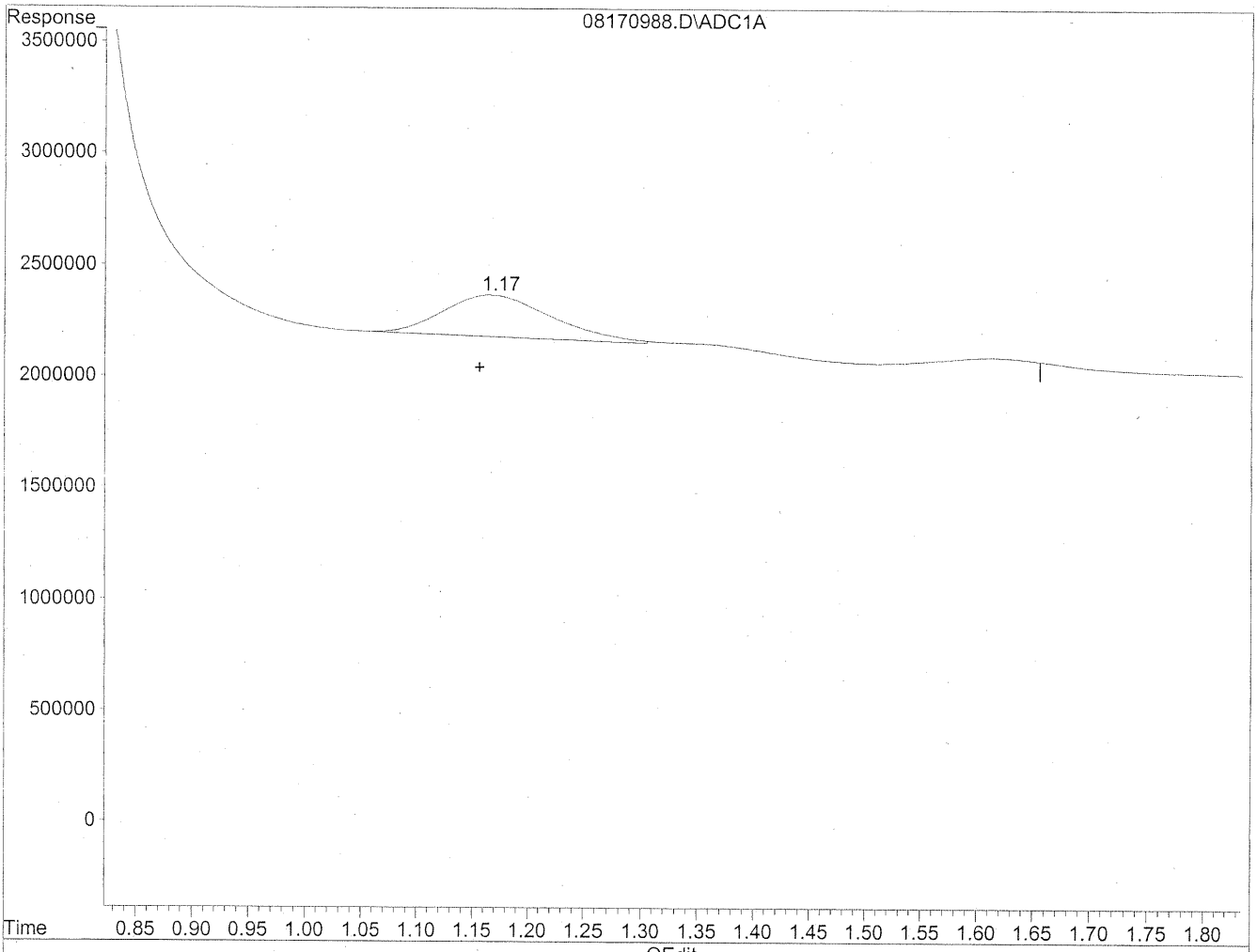
350

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170988.D Vial: 2
Acq On : 18 Aug 2009 12:38 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 13:16 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 68.626ng/ml m

response 12598459

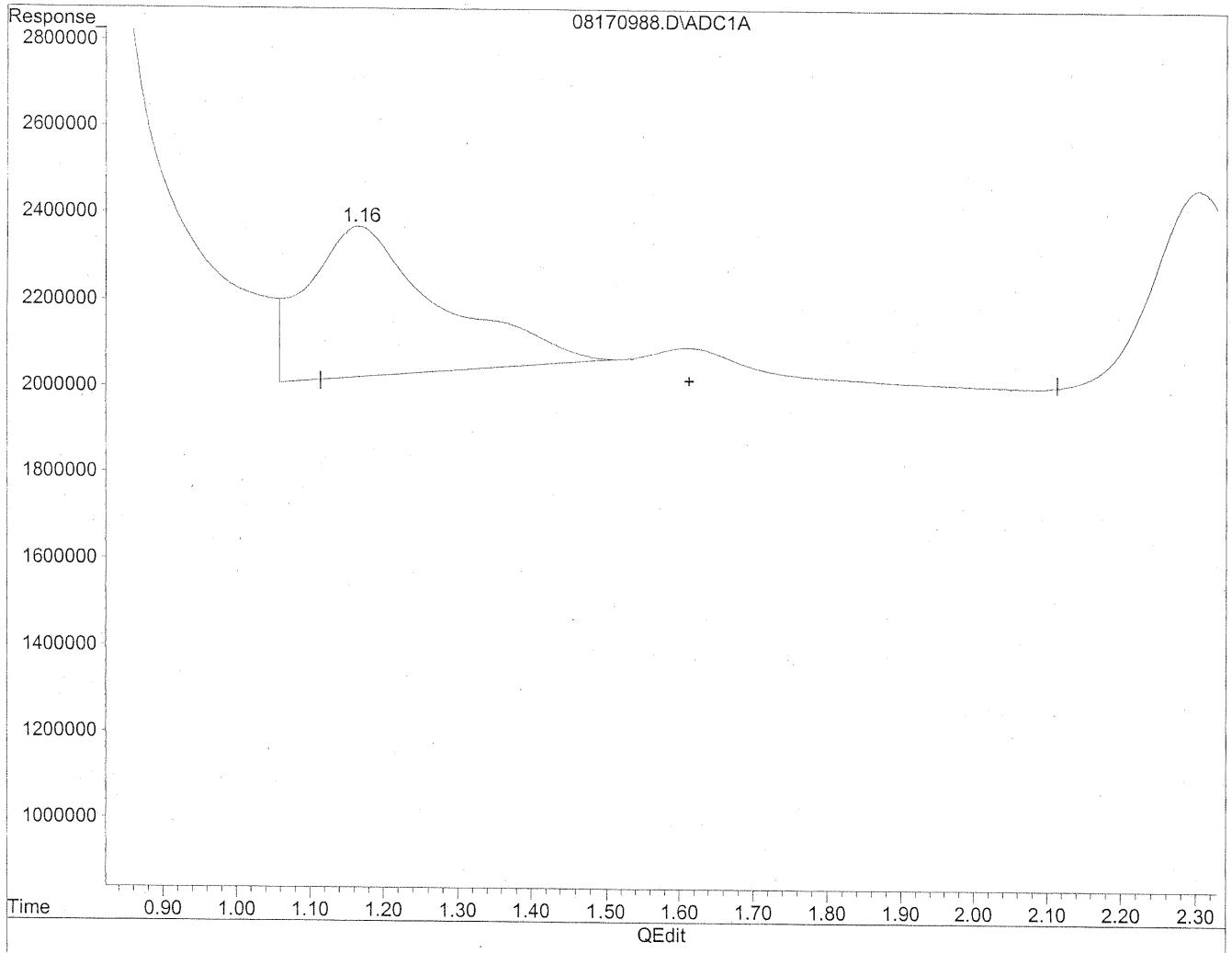
HC 8/22/09 LC KEB 8/23/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170988.D Vial: 2
Acq On : 18 Aug 2009 12:38 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 13:16 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.17min 308.521ng/ml

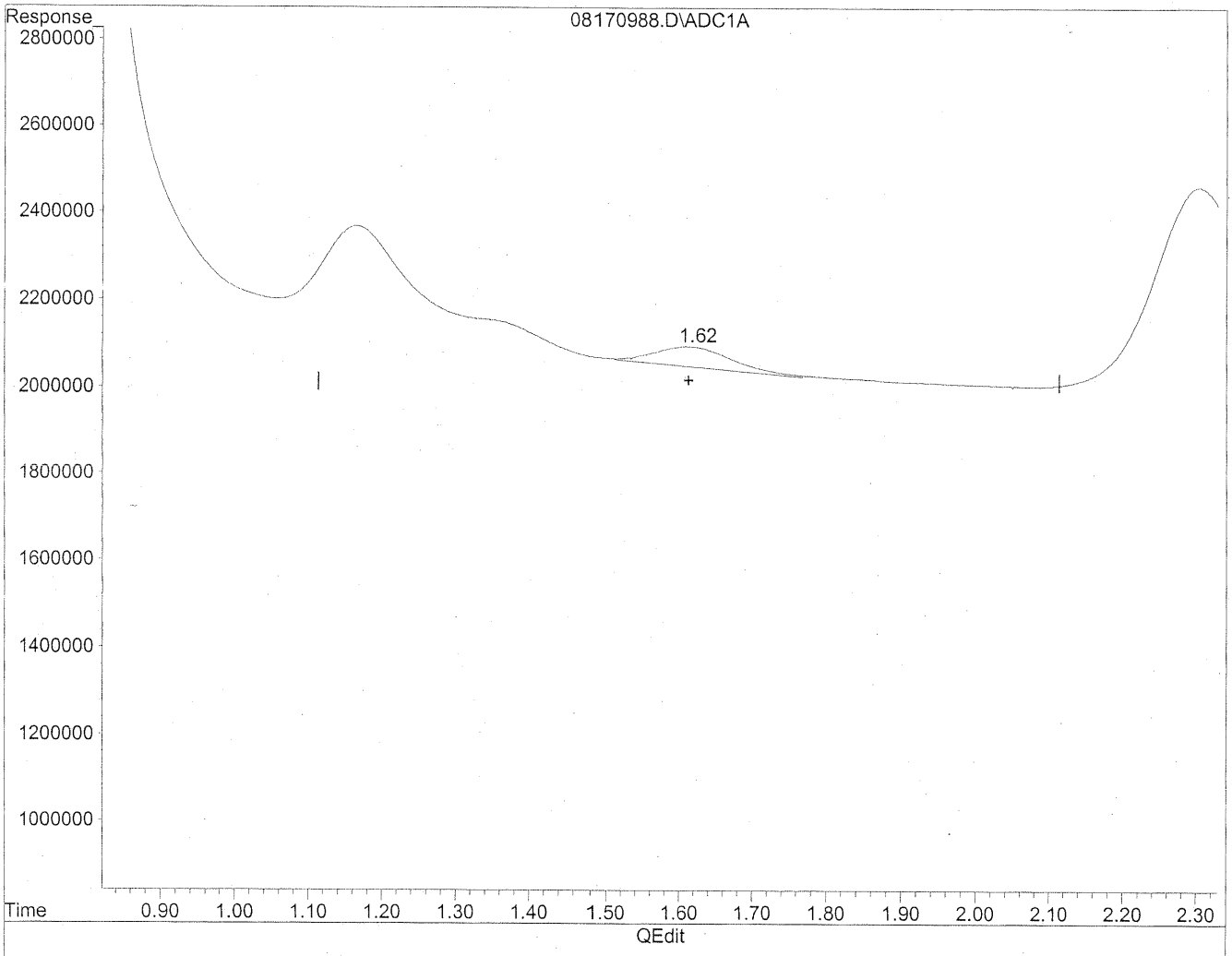
response 43261965

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170988.D Vial: 2
Acq On : 18 Aug 2009 12:38 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 13:16 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.62min 23.762ng/ml m

response 3332020

HC 8/22/07 mp

KB 8/26/07

353

(+) = Expected Retention Time

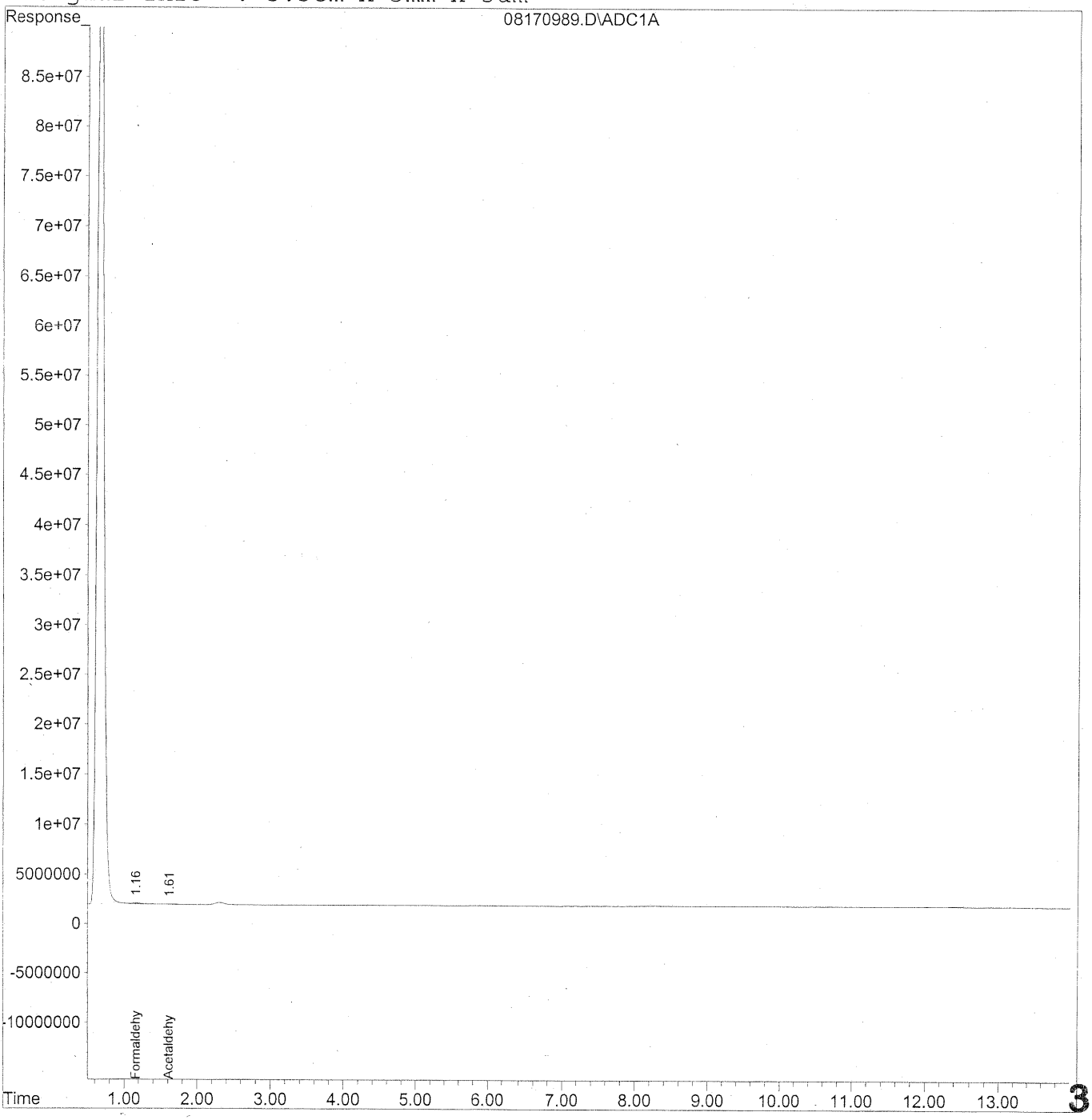
08170988.D TO110709.M Sat Aug 22 13:17:18 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170989.D Vial: 3
Acq On : 18 Aug 2009 12:53 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 13:17 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



354

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170989.D Vial: 3
Acq On : 18 Aug 2009 12:53 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 13:17 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

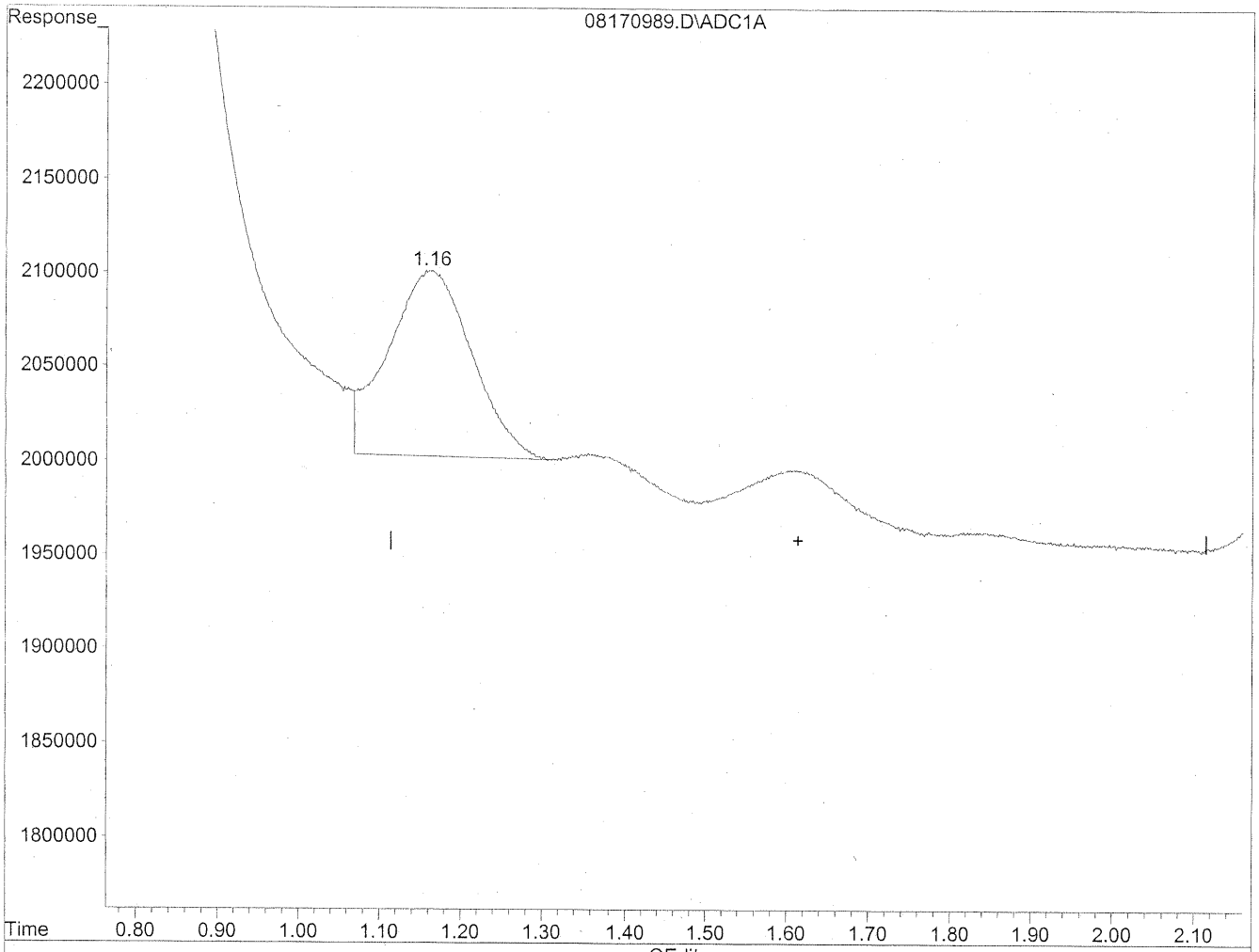
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	7058959	38.451	ng/ml
2)	Acetaldehyde	1.61	1866097	13.308	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170989.D Vial: 3
Acq On : 18 Aug 2009 12:53 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 13:17 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.16min 50.341ng/ml

response 7058959

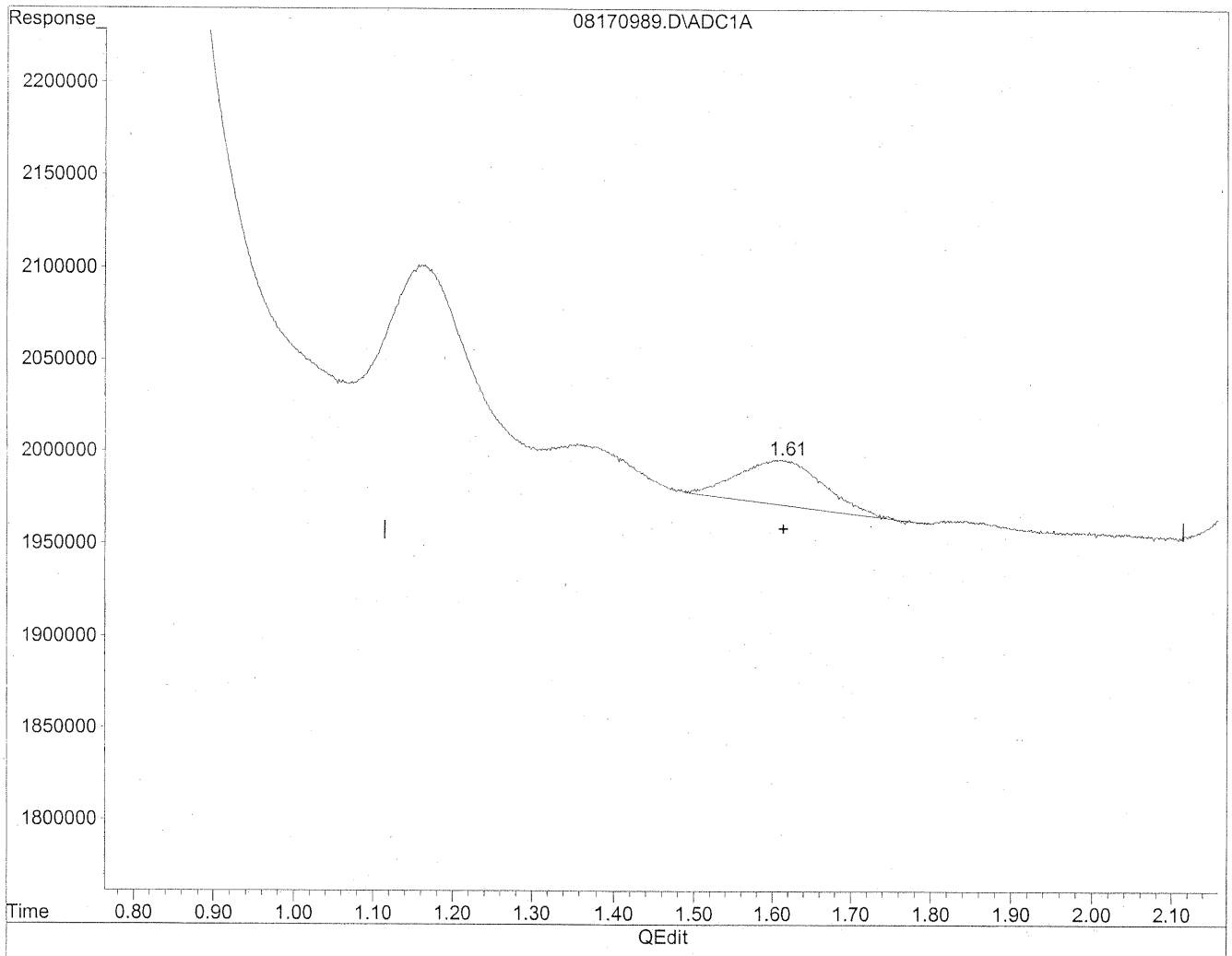
356

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170989.D Vial: 3
Acq On : 18 Aug 2009 12:53 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 13:17 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.61min 13.308ng/ml m

response 1866097

HC 8/22/09 MP KB 8/23/09

357

(+) = Expected Retention Time

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **Method Blank (19:24)**
Client Project ID: **16512**

CAS Project ID: P0902772
CAS Sample ID: P090818-MB

Test Code: EPA Method TO-11A
Instrument ID: Waters LC Module I Plus/UV_Vis 360/LC1
Analyst: Hani Cherazaie
Sampling Media: Silica Gel DNPH Tube
Test Notes: **BC**

Date Collected: NA
Date Received: NA
Date Analyzed: 08/18/09
Desorption Volume: 1.0 ml
Volume Sampled: NA Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	Formaldehyde	< 100	NA	NA	NA	NA	
75-07-0	Acetaldehyde	< 100	NA	NA	NA	NA	
123-38-6	Propionaldehyde	< 100	NA	NA	NA	NA	
4170-30-3	Crotonaldehyde, Total	< 100	NA	NA	NA	NA	
123-72-8	Butyraldehyde	< 100	NA	NA	NA	NA	
100-52-7	Benzaldehyde	< 100	NA	NA	NA	NA	
590-86-3	Isovaleraldehyde	< 100	NA	NA	NA	NA	
110-62-3	Valeraldehyde	< 100	NA	NA	NA	NA	
529-20-4	o-Tolualdehyde	< 100	NA	NA	NA	NA	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	NA	NA	NA	NA	
66-25-1	n-Hexaldehyde	< 100	NA	NA	NA	NA	
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	NA	NA	NA	NA	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
NA = Not applicable.

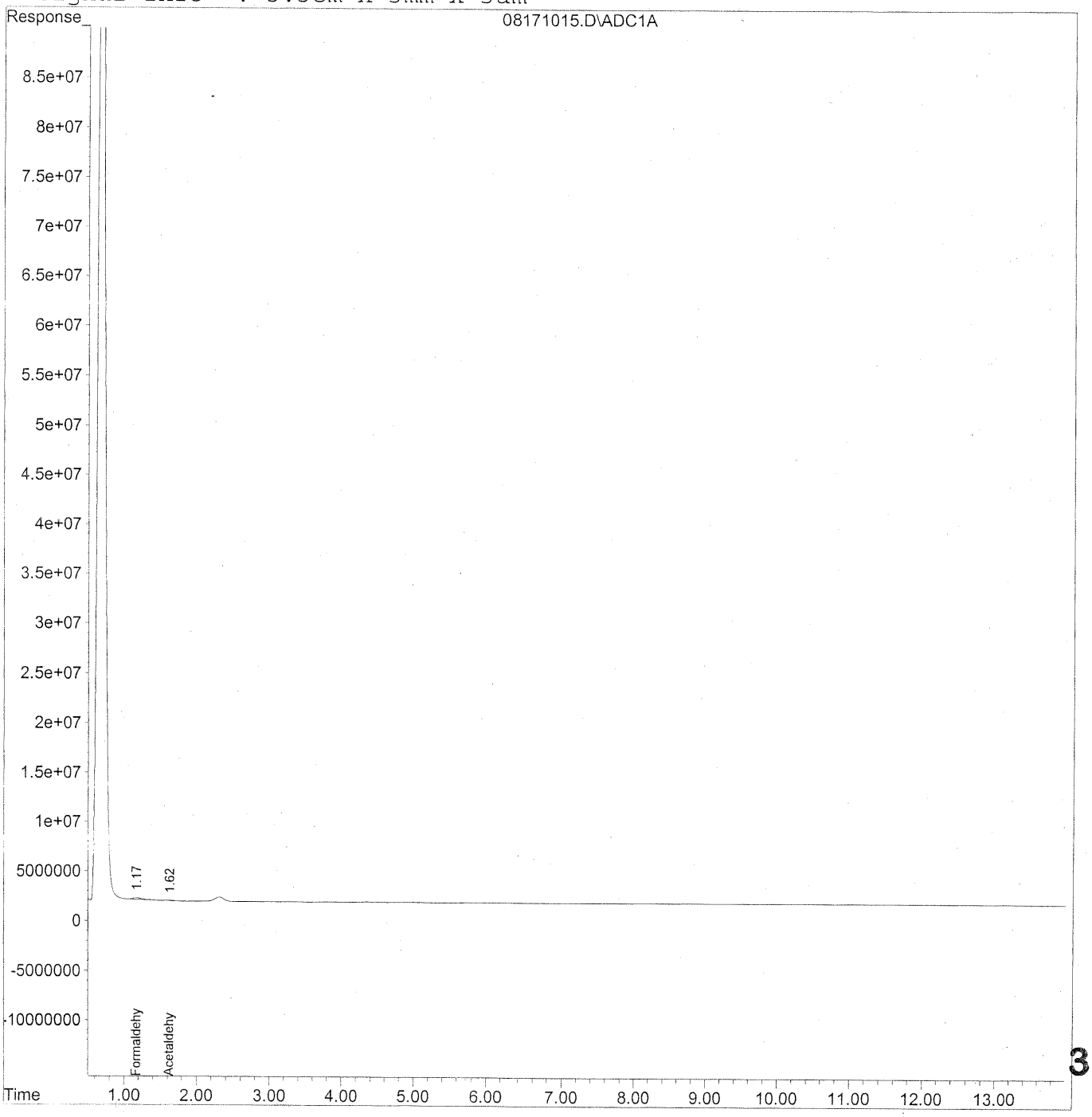
Verified By: _____ Date: 8/25

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171015.D Vial: 28
Acq On : 18 Aug 2009 7:24 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:30 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



359

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171015.D Vial: 28
Acq On : 18 Aug 2009 7:24 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:30 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

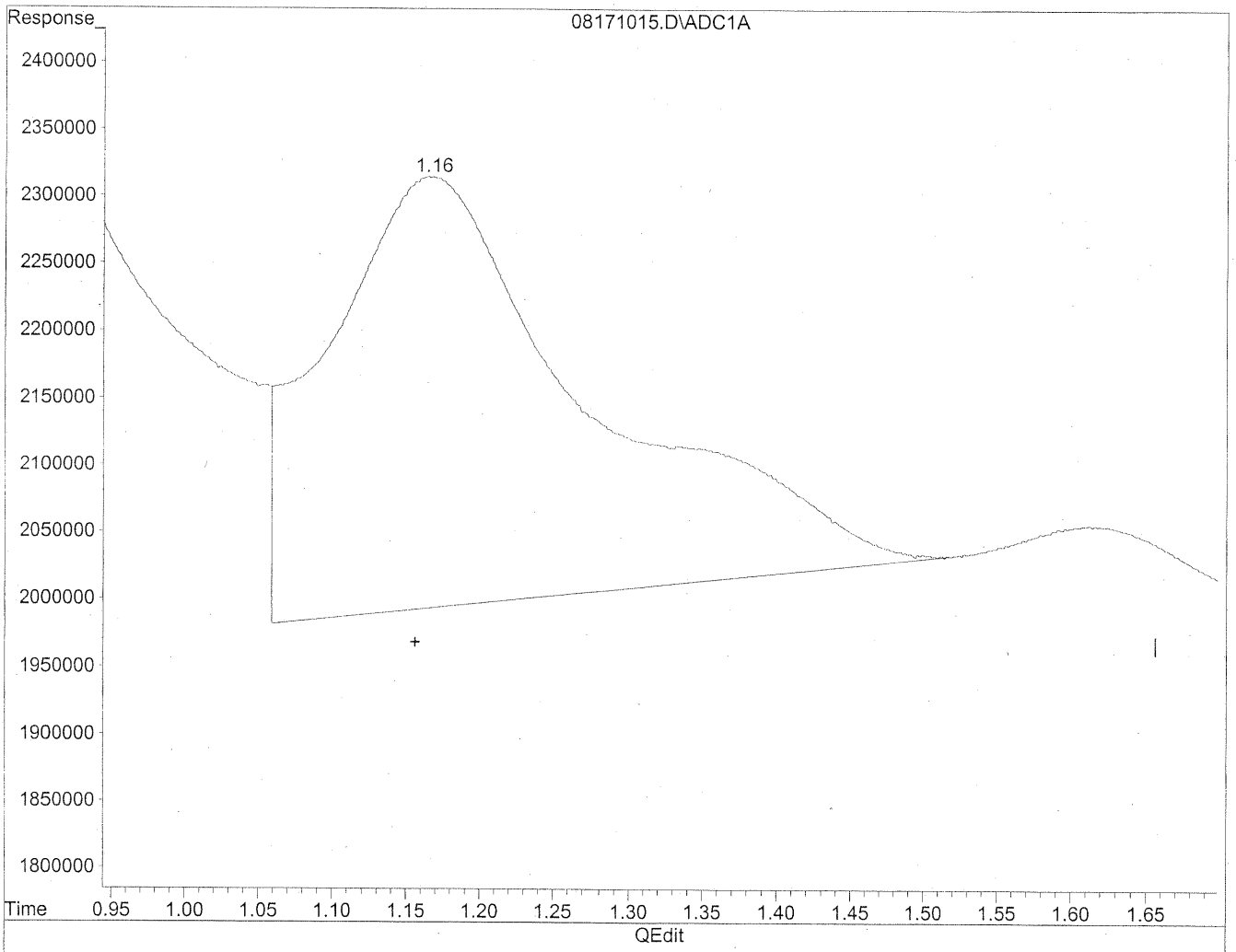
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	11331380	61.724	ng/mlm
2)	Acetaldehyde	1.62	2747549	19.594	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

(f)-RT Delta > 1/2 Window (m)

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171015.D Vial: 28
Acq On : 18 Aug 2009 7:24 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 216.252ng/ml

response 39699851

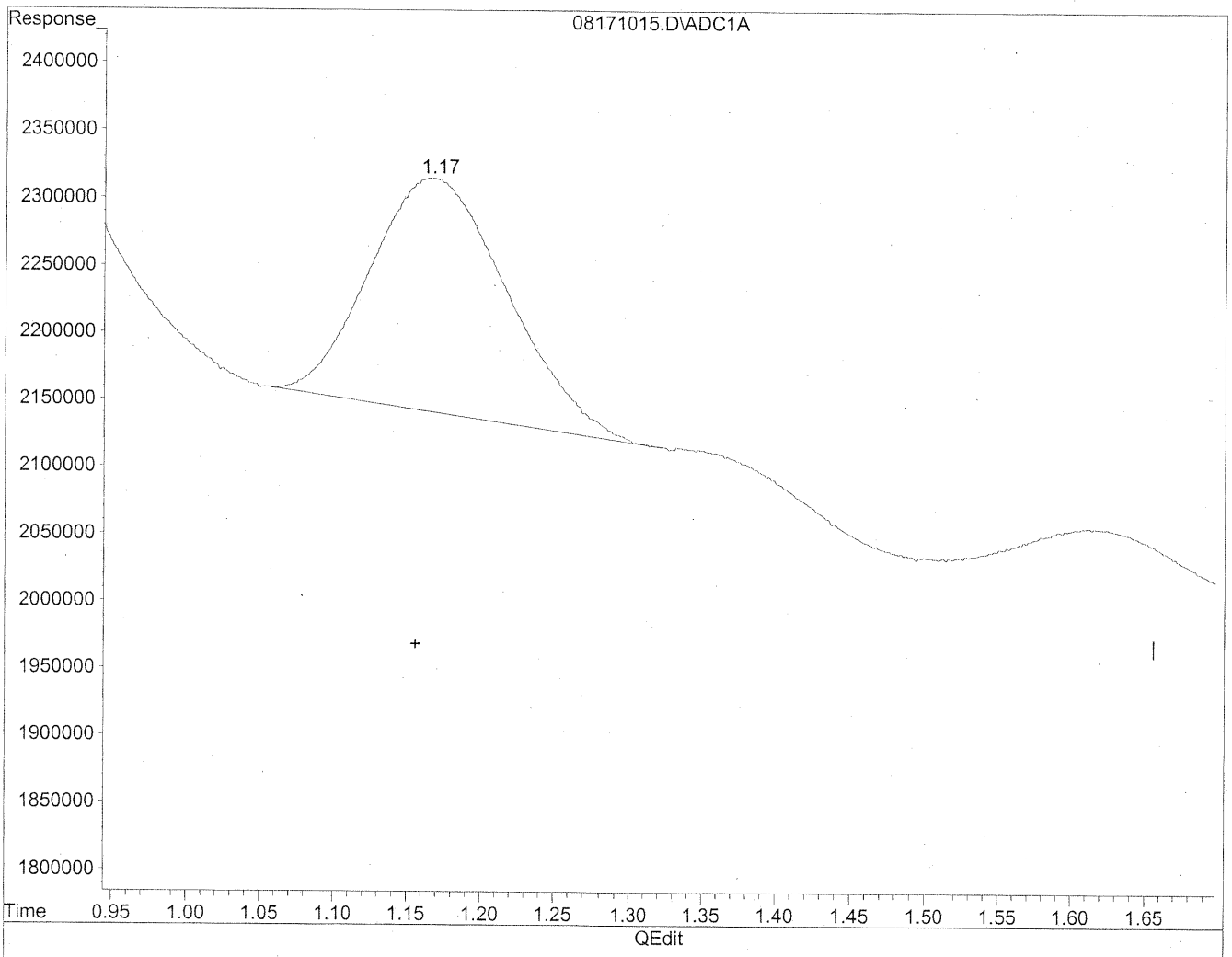
361

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171015.D Vial: 28
Acq On : 18 Aug 2009 7:24 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 61.724ng/ml m

response 11331380

tlc 8/22/09

lc 8/23/09

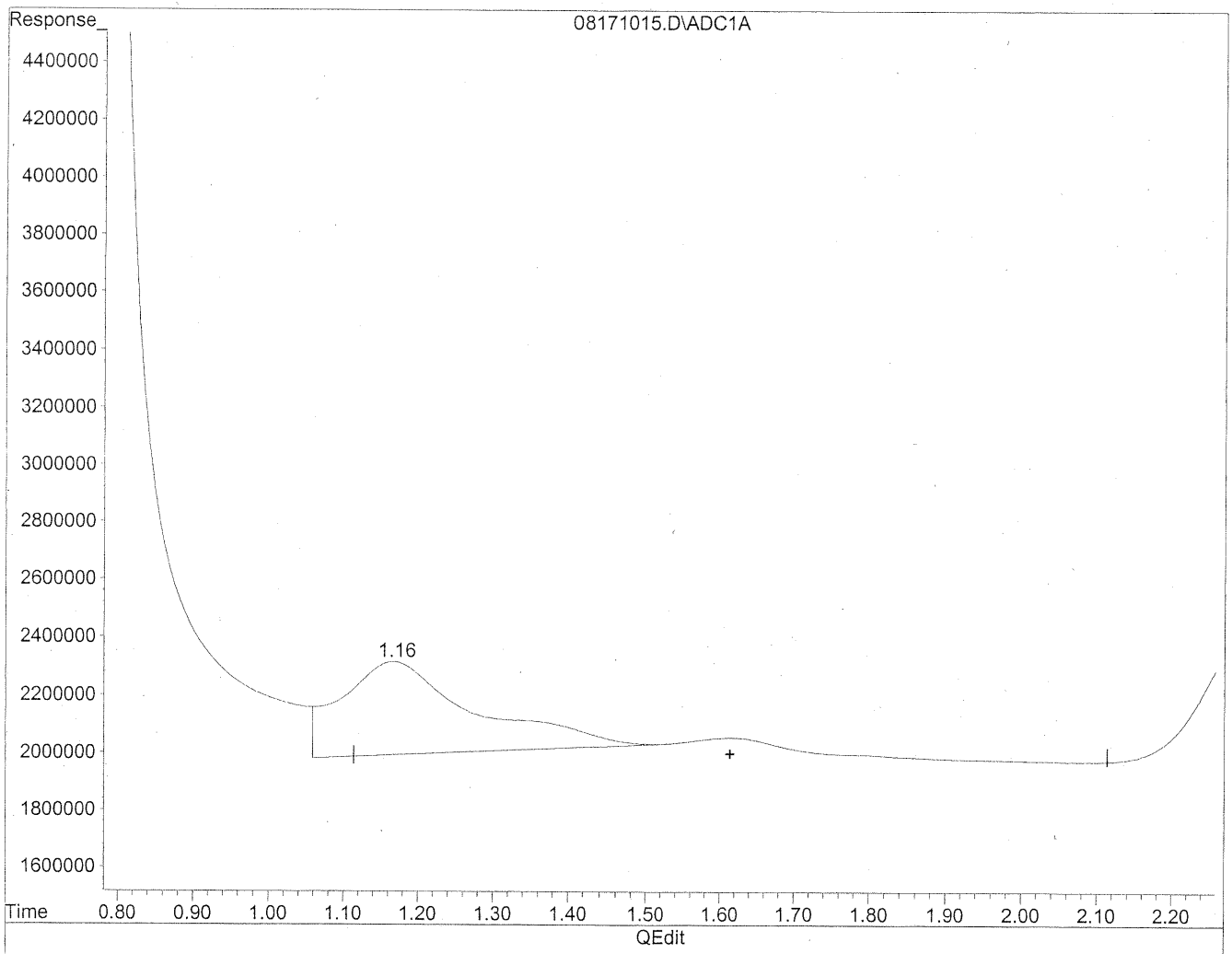
362

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171015.D Vial: 28
Acq On : 18 Aug 2009 7:24 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.17min 283.118ng/ml

response 39699851

363

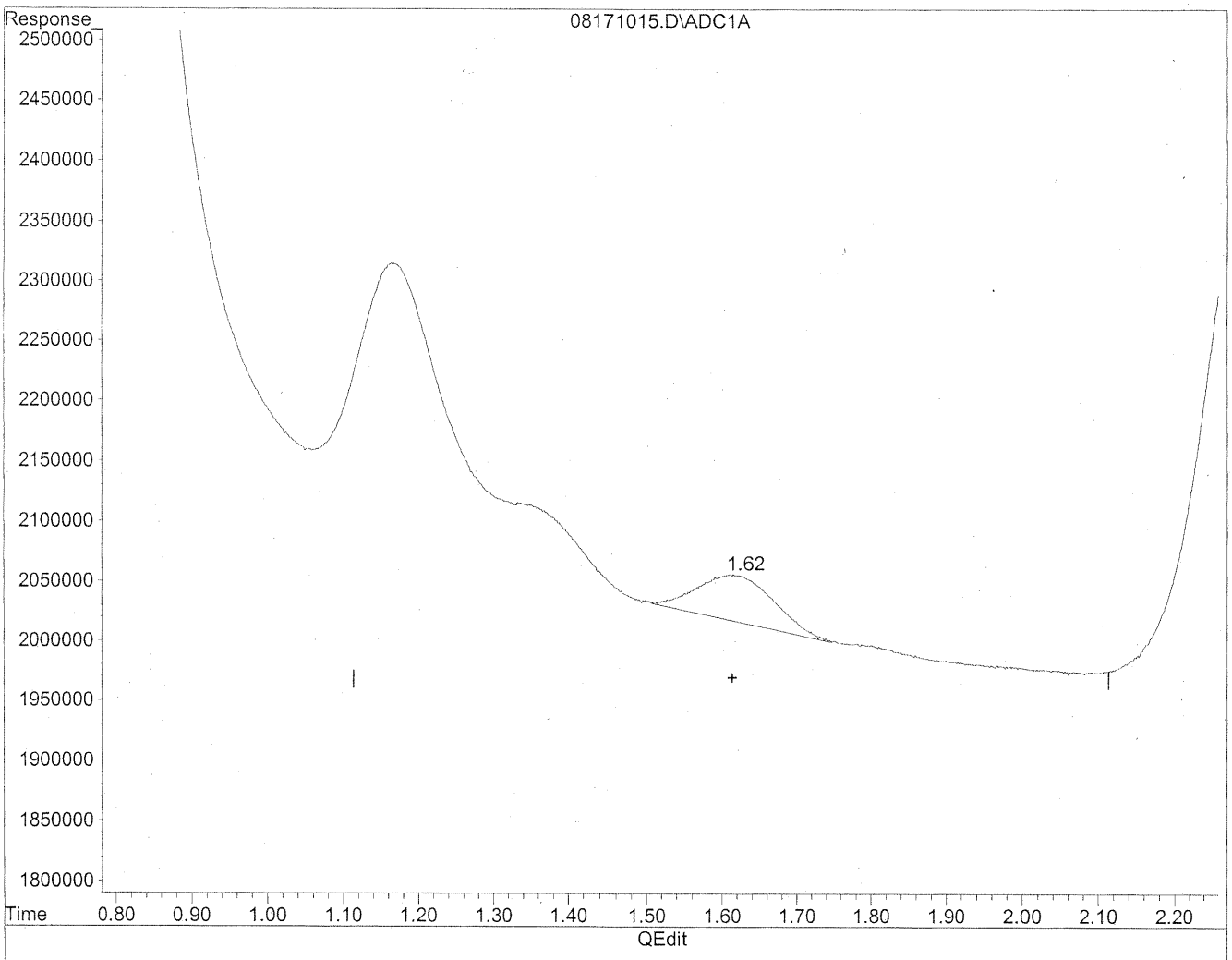
(+) = Expected Retention Time

08171015.D TO110709.M Sat Aug 22 14:30:31 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171015.D Vial: 28
Acq On : 18 Aug 2009 7:24 pm Operator: HC
Sample : MB front lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.62min 19.594ng/ml m

response 2747549

HC 8/22/09 MP

KL 8/23/09

364

(+) = Expected Retention Time

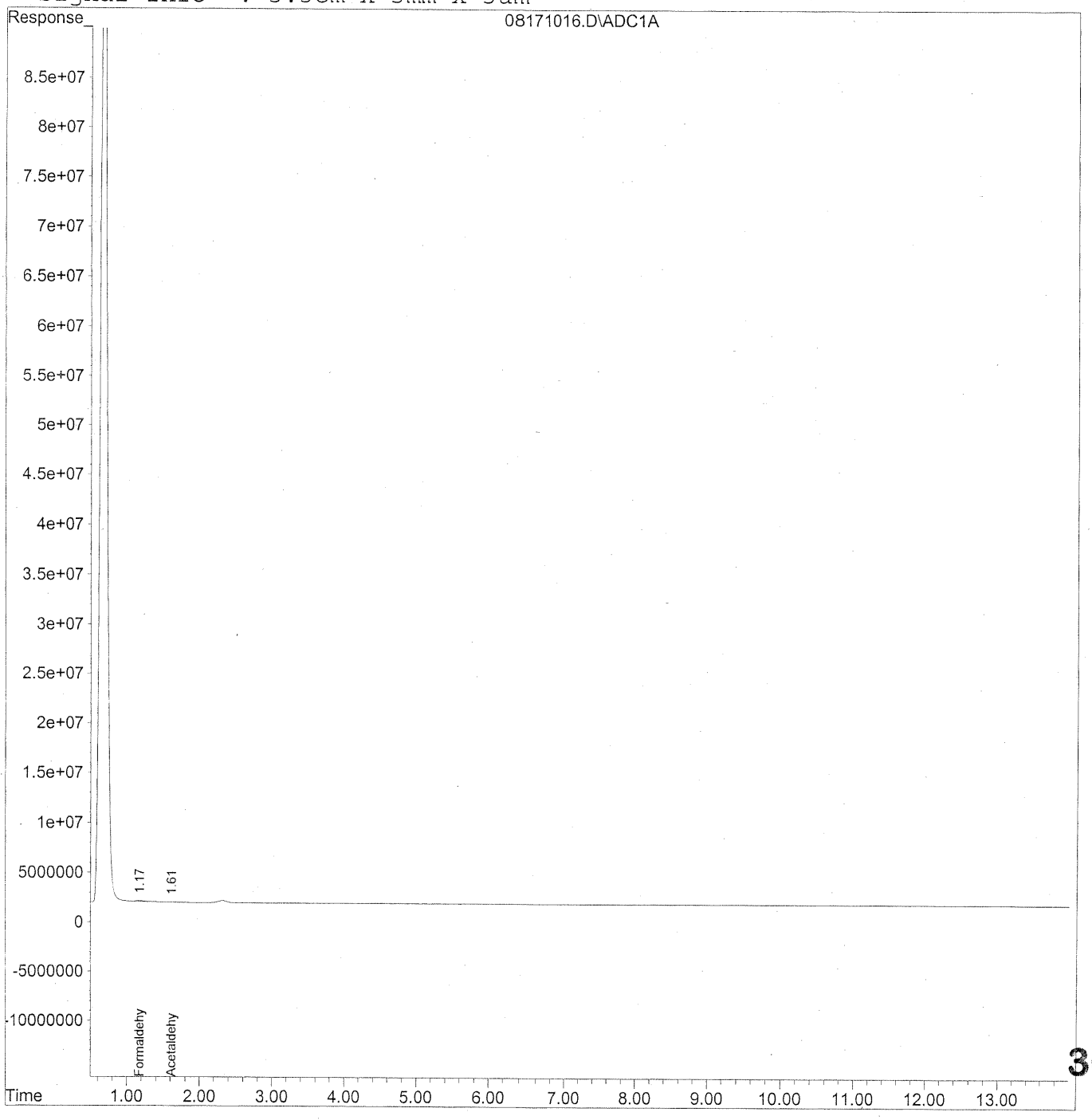
08171015.D TO110709.M Sat Aug 22 14:30:39 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171016.D Vial: 29
Acq On : 18 Aug 2009 7:39 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:31 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171016.D Vial: 29
Acq On : 18 Aug 2009 7:39 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 22 14:31 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

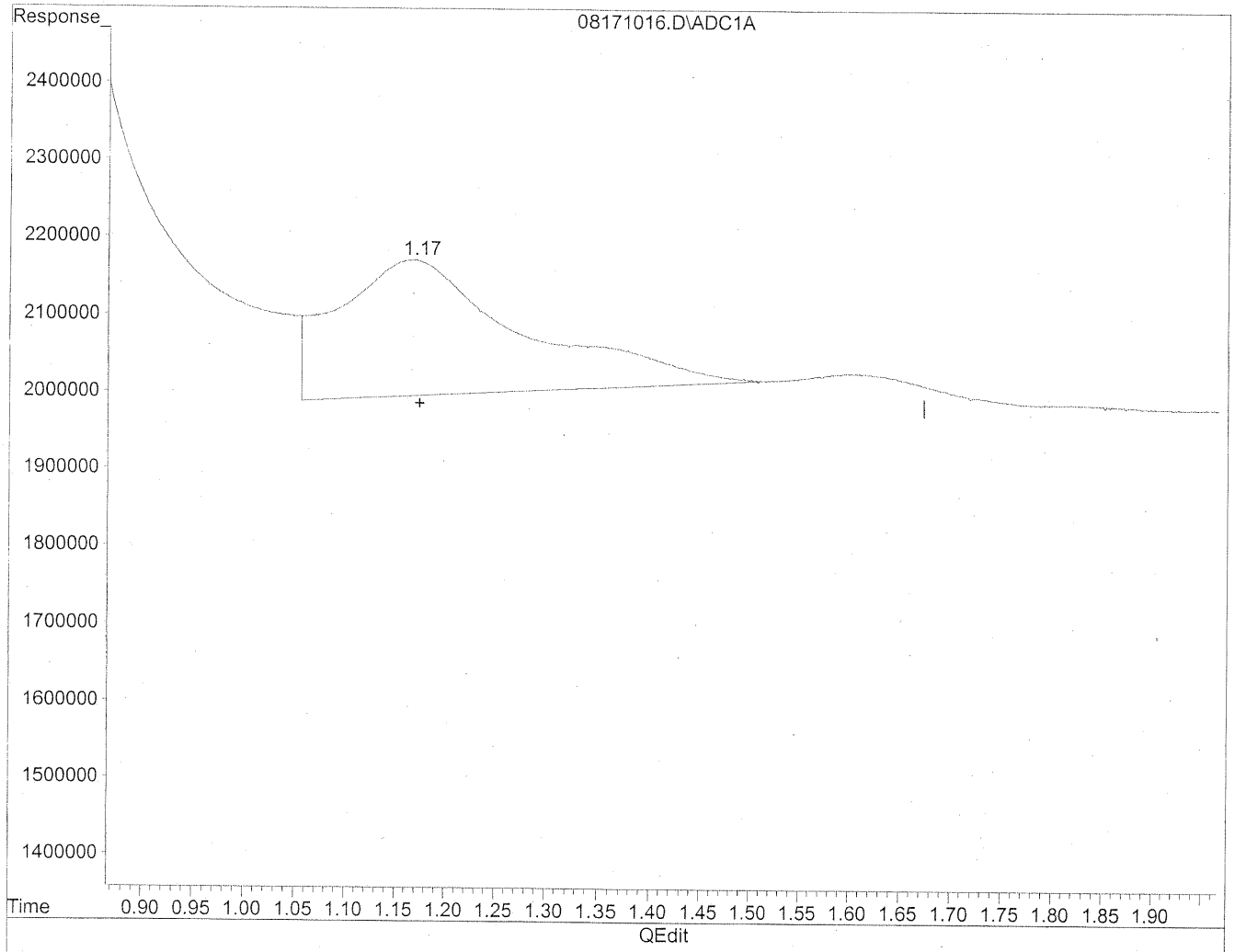
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	5788356	31.530	ng/mlm
2)	Acetaldehyde	1.61	1789412	12.761	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171016.D Vial: 29
Acq On : 18 Aug 2009 7:39 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 121.136ng/ml

response 22238308

367

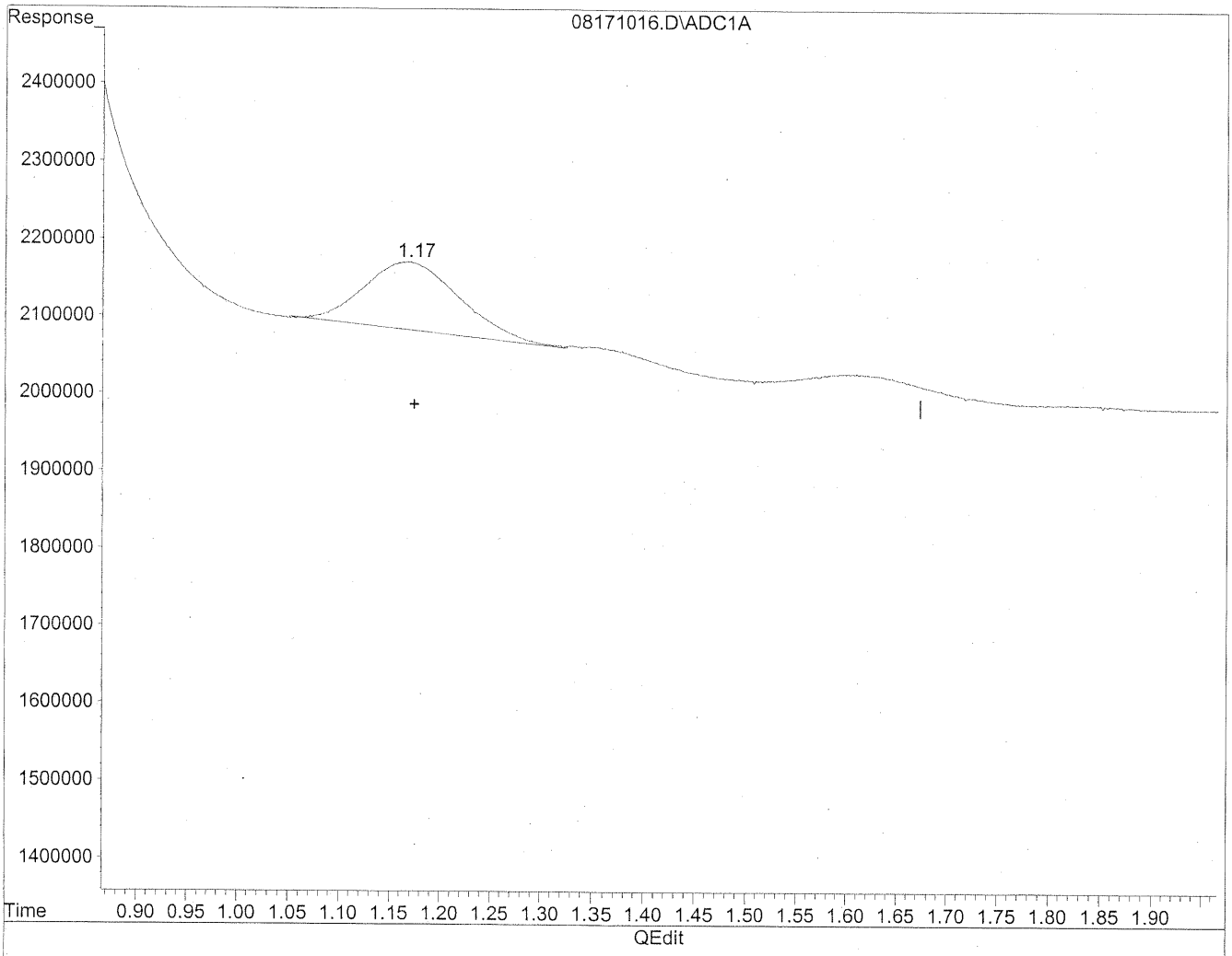
(+) = Expected Retention Time

08171016.D TO110709.M Sat Aug 22 14:30:51 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171016.D Vial: 29
Acq On : 18 Aug 2009 7:39 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.17min 31.530ng/ml m

response 5788356

HC 8/22/09 IC

KE 8/23/09

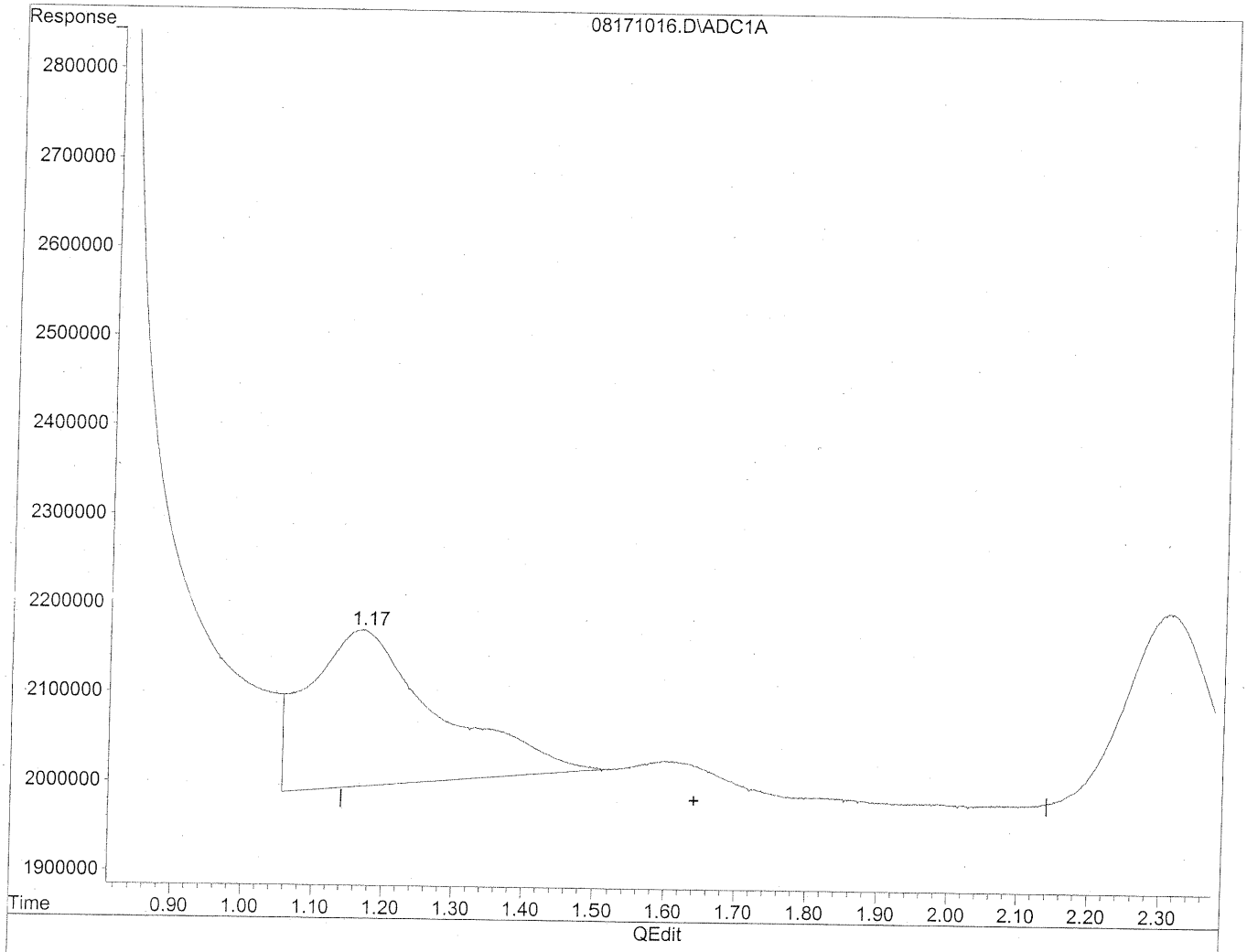
368

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171016.D Vial: 29
Acq On : 18 Aug 2009 7:39 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.17min 158.592ng/ml

response 22238308

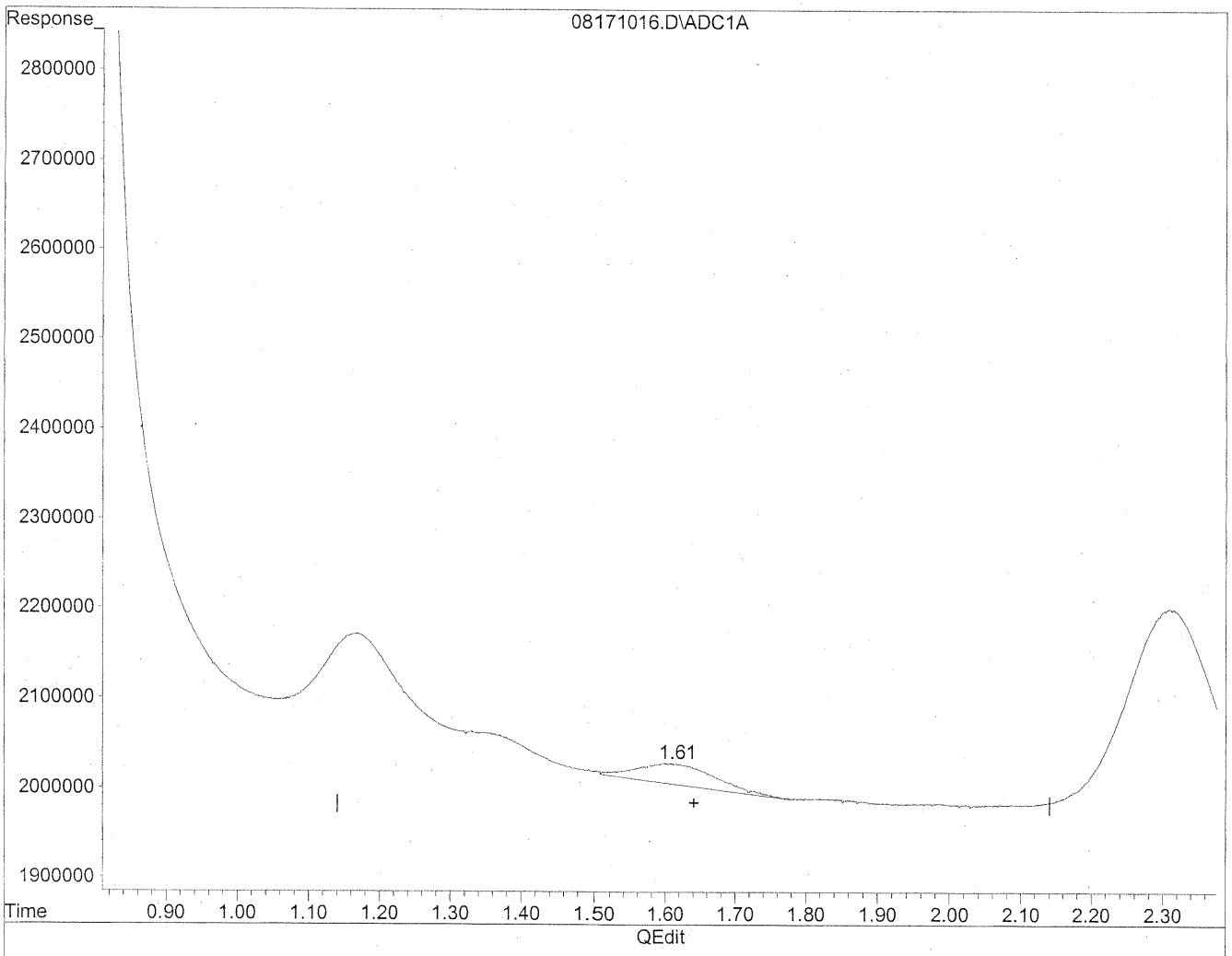
(+) = Expected Retention Time

08171016.D TO110709.M Sat Aug 22 14:21:02 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171016.D Vial: 29
Acq On : 18 Aug 2009 7:39 pm Operator: HC
Sample : MB back lot 6009/6097 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.61min 12.761ng/ml m

response 1789412

HC 8/22/09
MP
KE 8/28/09

370

(+) = Expected Retention Time

INITIAL CALIBRATION STANDARDS

Response Factor Report LC 01

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Mar 21 12:19:47 2005

Calibration Files
50 =07280905.D 100 =07280908.D 500 =07280909.D
1500 =07280912.D 5000 =02060917.D 10 =02060920.D

	Compound	50	100	500	1500	5000	10	Avg		%RSD
1)	Formaldehyde	1.776	1.838	1.825	1.831	1.848	1.897	1.836	E5	2.12
2)	Acetaldehyde	1.378	1.399	1.391	1.394	1.412	1.441	1.402	E5	1.55
3)	Propionaldehyde	1.021	1.096	1.057	1.058	1.074	1.096	1.067	E5	2.68
4)	Crotonaldehyde	1.082	0.953	0.945	0.944	0.951	0.969	0.974	E5	5.52
5)	Butyraldehyde	8.550	8.912	8.708	8.847	8.909	9.076	8.834	E4	2.07
6)	Benzaldehyde	6.116	6.908	6.719	6.549	6.563	6.666	6.587	E4	4.02
7)	Isovaleraldehyde	7.780	7.950	7.872	7.717	7.761	7.869	7.825	E4	1.11
8)	Valeraldehyde	7.609	7.695	7.248	7.114	7.160	7.276	7.351	E4	3.30
9)	o-Tolualdehyde	5.510	5.704	5.952	5.780	5.973	6.073	5.832	E4	3.55
10)	m,p-Tolualdehyde	5.048	5.565	5.415	5.370	5.457	5.541	5.400	E4	3.47
11)	Hexaldehyde	6.853	7.112	6.462	6.574	6.654	6.752	6.734	E4	3.41
12)	2,5-Dimethylbenzald	5.513	5.081	4.643	4.645	4.728	4.798	4.901	E4	6.95

HC 7/22/09

(#) = Out of Range

Calibration Status Report LC 01

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Jul 29 15:10:39 2009
Response via : Initial Calibration

#	ID	Conc	ISTD Conc	Path\File
1	50	50.00	0.00	J:\LC01\DATA\TO11\2009_07\28\07280905.D
2	100	100.00	0.00	J:\LC01\DATA\TO11\2009_07\28\07280908.D
3	500	500.00	0.00	J:\LC01\DATA\TO11\2009_07\28\07280909.D
4	1500	1500.00	0.00	J:\LC01\DATA\TO11\2009_07\28\07280912.D
5	5000	5000.00	0.00	J:\LC01\DATA\TO11\2009_07\28\07280915.D
6	10	10000.00	0.00	J:\LC01\DATA\TO11\2009_07\28\07280918.D

#	ID	Update Time	Quant Time	Acquisition Time
1	50	Jul 28 10:27 2009	Jul 28 10:27 19109	28 Jul 2009 9:39 am
2	100	Jul 28 14:52 2009	Jul 28 14:34 19109	28 Jul 2009 10:24 am
3	500	Jul 28 14:52 2009	Jul 28 14:40 19109	28 Jul 2009 10:39 am
4	1500	Jul 28 17:22 2009	Jul 28 14:45 19109	28 Jul 2009 11:24 am
5	5000	Jul 29 15:10 2009	Jul 28 14:48 19109	28 Jul 2009 12:09 pm
6	10	Jul 29 15:10 2009	Jul 28 14:49 19109	28 Jul 2009 12:54 pm

TO110709.M Wed Jul 29 15:10:44 2009

COLUMBIA ANALYTICAL SERVICES, INC.

Method: TO-11A
Analyst: HC

Printed : 7/30/09
Instrument : LC#1
Date Analysis : 6/23/00
Detector : UV-VIS 360
Sample Amount : 5ul

CALIBRATION RESPONSE FACTOR SUMMARY

Calibration Level	Form-Aldehyde	% rpd	Acet-Aldehyde	% rpd	Propion-Aldehyde	% rpd	Croton-Aldehyde	% rpd	Butyr-Aldehyde	% rpd	Benz-Aldehyde	% rpd
50ng/ml TO11A Std	8477013	4.54%	6307171	8.47%	4892636	4.12%	5507079	1.75%	4412295	3.21%	3362429	9.96%
50ng/ml TO11A Std	8859457	0.24%	6975740	1.23%	4973947	2.53%	4974991	8.08%	4293221	0.43%	3079204	0.70%
50ng/ml TO11A Std	9305088	4.78%	7389770	7.24%	5442713	6.66%	5754474	6.32%	4119144	3.64%	2732056	10.66%
100ng/ml TO11A St	18283557	0.51%	13784712	1.44%	10870707	0.86%	9346475	1.91%	8839595	0.81%	7282249	5.41%
100ng/ml TO11A St	18449443	0.39%	14434553	3.21%	11389784	3.88%	9814490	3.00%	9432197	5.84%	6706722	2.92%
100ng/ml TO11A St	18400032	0.12%	13737532	1.77%	10633406	3.02%	9424529	1.09%	8463028	5.03%	6735919	2.50%
500ng/ml TO11A St	91593554	0.39%	70468869	0.90%	53468174	1.20%	47866960	1.26%	43271557	0.62%	32616313	2.91%
500ng/ml TO11A St	90711575	0.57%	69140255	1.00%	52850412	0.03%	47584179	0.66%	43677338	0.31%	34085310	1.46%
500ng/ml TO11A St	91399555	0.18%	69908753	0.10%	52190620	1.22%	46362546	1.92%	43673214	0.30%	34084716	1.46%
1500ng/ml TO11A S	275380897	0.26%	209374751	0.16%	159030091	0.21%	143227783	1.11%	134132687	1.08%	9887868	0.65%
1500ng/ml TO11A S	274724982	0.02%	209301649	0.12%	158919579	0.14%	142112419	0.32%	132549734	0.12%	9818365	0.06%
1500ng/ml TO11A S	273895978	0.28%	208465321	0.28%	158125683	0.36%	139629551	1.43%	131425702	0.96%	9765243	0.60%
5000ng/ml TO11A S	928364658	0.45%	706170560	0.05%	539067854	0.39%	476268543	0.19%	446392739	0.21%	328286106	0.04%
5000ng/ml TO11A S	925768000	0.17%	708552415	0.38%	540133923	0.59%	477844499	0.52%	446568052	0.25%	328413551	0.08%
5000ng/ml TO11A S	918424042	0.62%	702791887	0.43%	535167082	0.98%	471954575	0.72%	443441833	0.45%	327762901	0.12%
10000ng/ml TO11A	1908653125	0.62%	1450154617	0.67%	1099941045	0.36%	972691462	0.37%	910896701	0.36%	668462127	0.28%
10000ng/ml TO11A	1905913073	0.48%	1446499891	0.41%	1098837646	0.26%	971357788	0.23%	911328243	0.41%	669128969	0.38%
10000ng/ml TO11A	1875917434	1.10%	1425028469	1.08%	1089338811	0.61%	963283335	0.60%	900561239	0.78%	662238443	0.66%

AVERAGE RESPONSE FACTOR

Method:
Analyst:

CALIBRATION F

Calibration Level	Isovaler-Aldehyde	% rpd	Valer-Aldehyde	% rpd	o-Tolu-Aldehyde	% rpd	m,p-Tolu-Aldehyde	% rpd	Hex-Aldehyde	% rpd	2,5-Dimethl Benz-Aldehyde	% rpd
50ng/ml TO11A Std	4167653	7.13%	3532734	7.15%	338/183	22.94%	5445142	7.87%	3244418	5.31%	2546144	7.64%
50ng/ml TO11A Std	4002738	2.89%	4025564	5.81%	2461625	10.65%	489/087	2.98%	3295067	3.83%	2605446	5.49%
50ng/ml TO11A Std	3500271	10.02%	3855749	1.34%	2416389	12.29%	4801019	4.89%	3739368	9.14%	3118537	13.13%
100ng/ml TO11A St	7487274	5.83%	7060988	8.24%	5548699	2.73%	109/9457	1.36%	6702769	5.76%	5798505	14.13%
100ng/ml TO11A St	8338385	4.88%	811/341	5.49%	5921917	3.82%	11235135	0.94%	7714022	8.46%	4735227	6.80%
100ng/ml TO11A St	8025579	0.95%	7906862	2.75%	5642221	1.09%	1117/259	0.42%	6920120	2.70%	4707951	7.33%
500ng/ml TO11A St	37944016	3.60%	3557/4509	1.84%	29317615	1.49%	5327/4975	1.62%	32888440	1.80%	23823948	2.62%
500ng/ml TO11A St	40968120	4.08%	36648075	1.12%	29793454	0.11%	54514161	0.67%	31855201	1.40%	22510750	3.03%
500ng/ml TO11A St	39175205	0.48%	36501988	0.72%	30169058	1.37%	54668231	0.95%	32179520	0.40%	23309464	0.41%
1500ng/ml TO11A S	115866442	0.09%	107104204	0.36%	86339652	0.42%	162946532	1.14%	98895406	0.29%	69932636	0.37%
1500ng/ml TO11A S	116723586	0.83%	107107592	0.37%	85940120	0.88%	161094009	0.01%	98090122	0.53%	6887/3541	1.15%
1500ng/ml TO11A S	114690000	0.92%	10593/177	0.73%	8782422/	1.30%	159292531	1.13%	98846718	0.24%	70224395	0.79%
5000ng/ml TO11A S	38824/386	0.05%	357832844	0.04%	298513860	0.05%	545640330	0.02%	332315493	0.11%	235692401	0.30%
5000ng/ml TO11A S	388941560	0.23%	35967/6615	0.47%	30007/384	0.48%	547211501	0.27%	333701808	0.31%	237108293	0.30%
5000ng/ml TO11A S	386992833	0.28%	356464469	0.43%	29737/4461	0.43%	544331756	0.26%	332038452	0.19%	236428207	0.01%
10000ng/ml TO11A	79032817	0.44%	7302186/3	0.36%	60820827/6	0.16%	1111180147	0.26%	67351680/	0.25%	47846094/	0.27%
10000ng/ml TO11A	78802619/0	0.15%	72983921/0	0.31%	61032623/8	0.50%	1113209810	0.45%	68191578/5	0.99%	4847/6391/8	1.04%
10000ng/ml TO11A	78225680/4	0.59%	7227/49626	0.67%	60325659/9	0.66%	11003845/3	0.71%	67019336/0	0.74%	47/6113656	0.76%

AVERAGE RESI

	Form-Aldehyde	Acet-Aldehyde	Propion-Aldehyde	Croton-Aldehyde	Butyr-Aldehyde	Benz-Aldehyde
50ng/ml TO11A St	8880519	6890894	5103099	5412181	4274887	3057896
100ng/ml TO11A S	18377677	13985599	10964632	9528498	8911607	6908297
500ng/ml TO11A S	91234895	69839292	52836402	47271228	43540703	33595446
1500ng/ml TO11A	274667286	209047240	158691784	141656584	132702708	98238389
5000ng/ml TO11A	924185567	705838287	536958953	475355872	445467541	328154186
10000ng/ml TO11A	1896827877	1440560992	1096039167	969110862	907595394	666609846

	Isovaler-Aldehyde	Valer-Aldehyde	o-Tolu-Aldehyde	m,p-Tolu-Aldehyde	Hex-Aldehyde	2,5-Dimethl benz-Aldehyde
50ng/ml TO11A St	3890221	3804682	2755066	5047749	3426284	2756709
100ng/ml TO11A S	7950413	7695064	5704279	11130617	7112304	5080561
500ng/ml TO11A S	39362447	36241524	29760042	54152456	32307720	23214721
1500ng/ml TO11A	115760009	106716324	86701333	161111024	98610749	69676857
5000ng/ml TO11A	388060593	357991309	298655235	545727862	332685251	236409634
10000ng/ml TO11A	786870437	727602503	607263704	1108258177	675208651	479779507

TO-11A CALIBRATION STANDARDS LIST							
50ng/ml TO11A Std S21-07270908							
100ng/ml TO11A Std S21-07270905							
500ng/ml TO11A Std S21-07270904							
1500ng/ml TO11A Std S21-07270903							
5000ng/ml TO11A Std S21-07270902							
10000ng/ml TO11A Std S21-07270901							

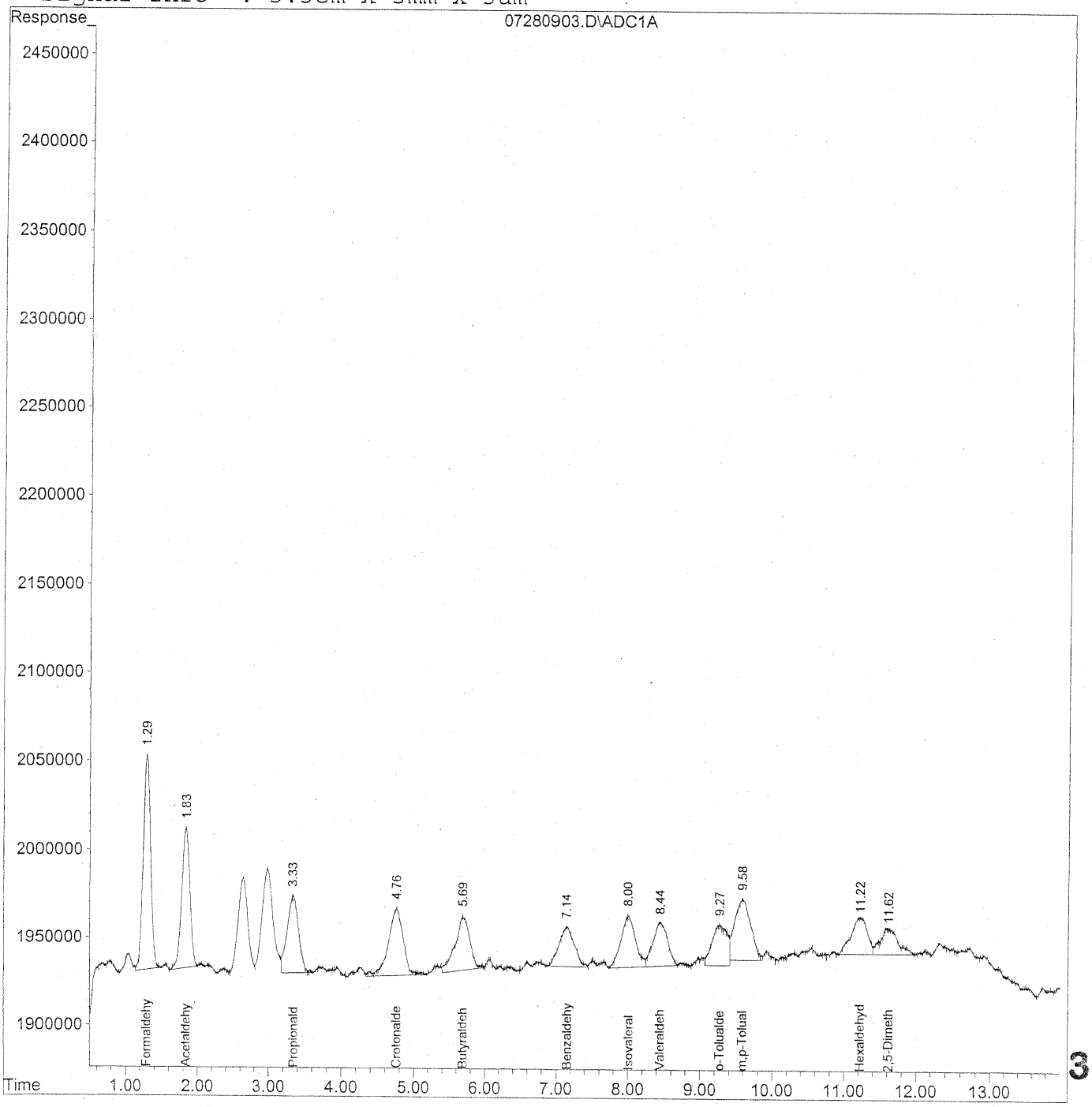
HC 4/29/05

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280903.D Vial: 3
Acq On : 28 Jul 2009 9:09 am Operator: HC
Sample : 50ng/ml TO11A Std S21-07270908 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:22 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



379

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280903.D Vial: 3
Acq On : 28 Jul 2009 9:09 am Operator: HC
Sample : 50ng/ml TO11A Std S21-07270908 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:22 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

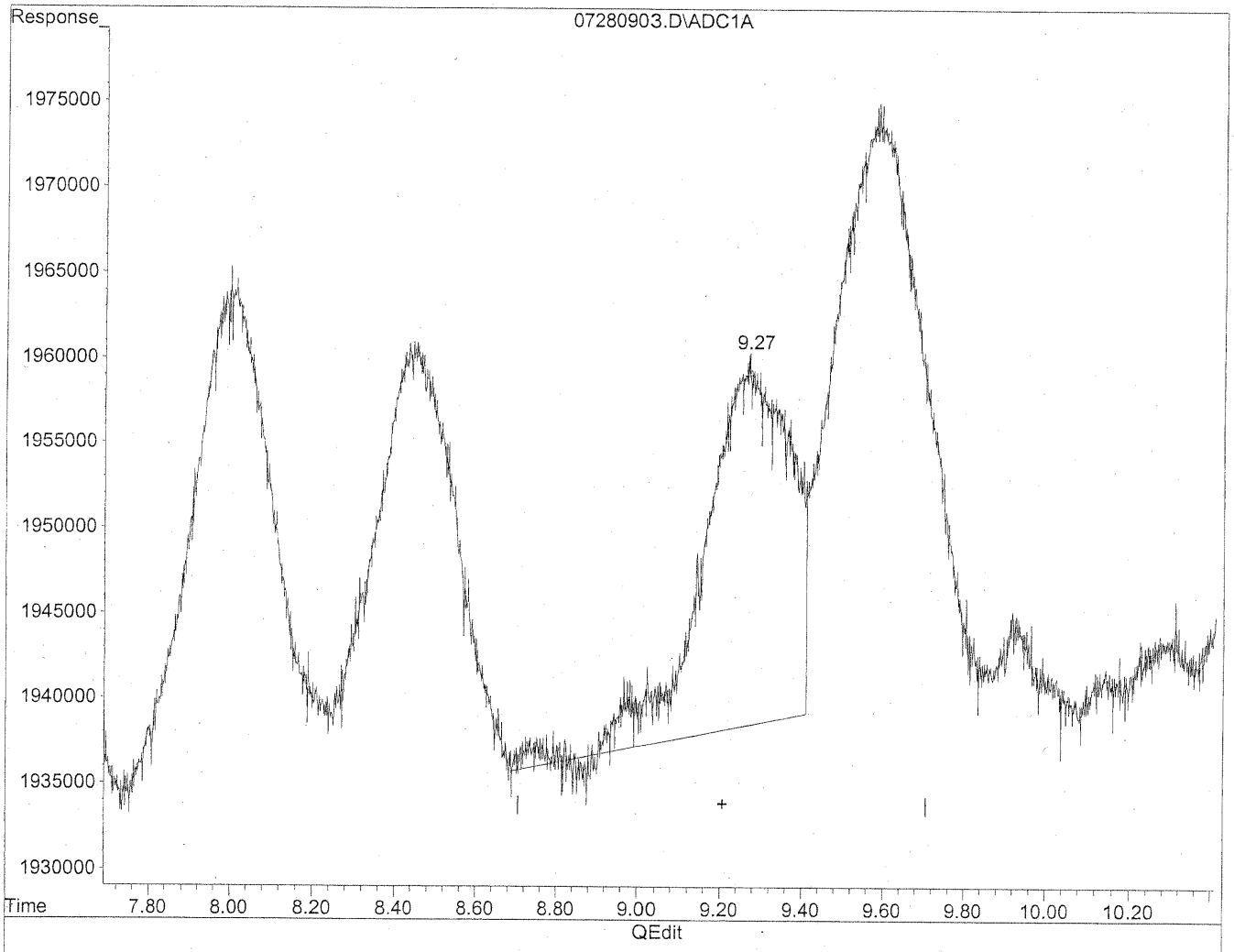
	Compound	R.T.	Response	Conc Units
	Target Compounds			
1)	Formaldehyde	1.29	8477013	48.277 ng/ml
2)	Acetaldehyde	1.83	6307171	46.755 ng/ml
3)	Propionaldehyde	3.34	4892636	47.596 ng/ml
4)	Crotonaldehyde	4.76	5507079	49.813 ng/ml
5)	Butyraldehyde	5.70	4412295	54.828 ng/ml
6)	Benzaldehyde	7.15	3362429	53.310 ng/ml
7)	Isovaleraldehyde	8.01	4167653	47.012 ng/ml
8)	Valeraldehyde	8.45	3532734	42.514 ng/ml
9)	o-Tolualdehyde	9.27	3387183	62.877 ng/mlm
10)	m,p-Tolualdehyde	9.58	5445142	101.089 ng/mlm
11)	Hexaldehyde	11.22	3244418	48.324 ng/ml
12)	2,5-Dimethylbenzaldehyde	11.61	2546144	49.027 ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280903.D Vial: 3
Acq On : 28 Jul 2009 9:09 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.27min 57.721ng/ml

response 3109441

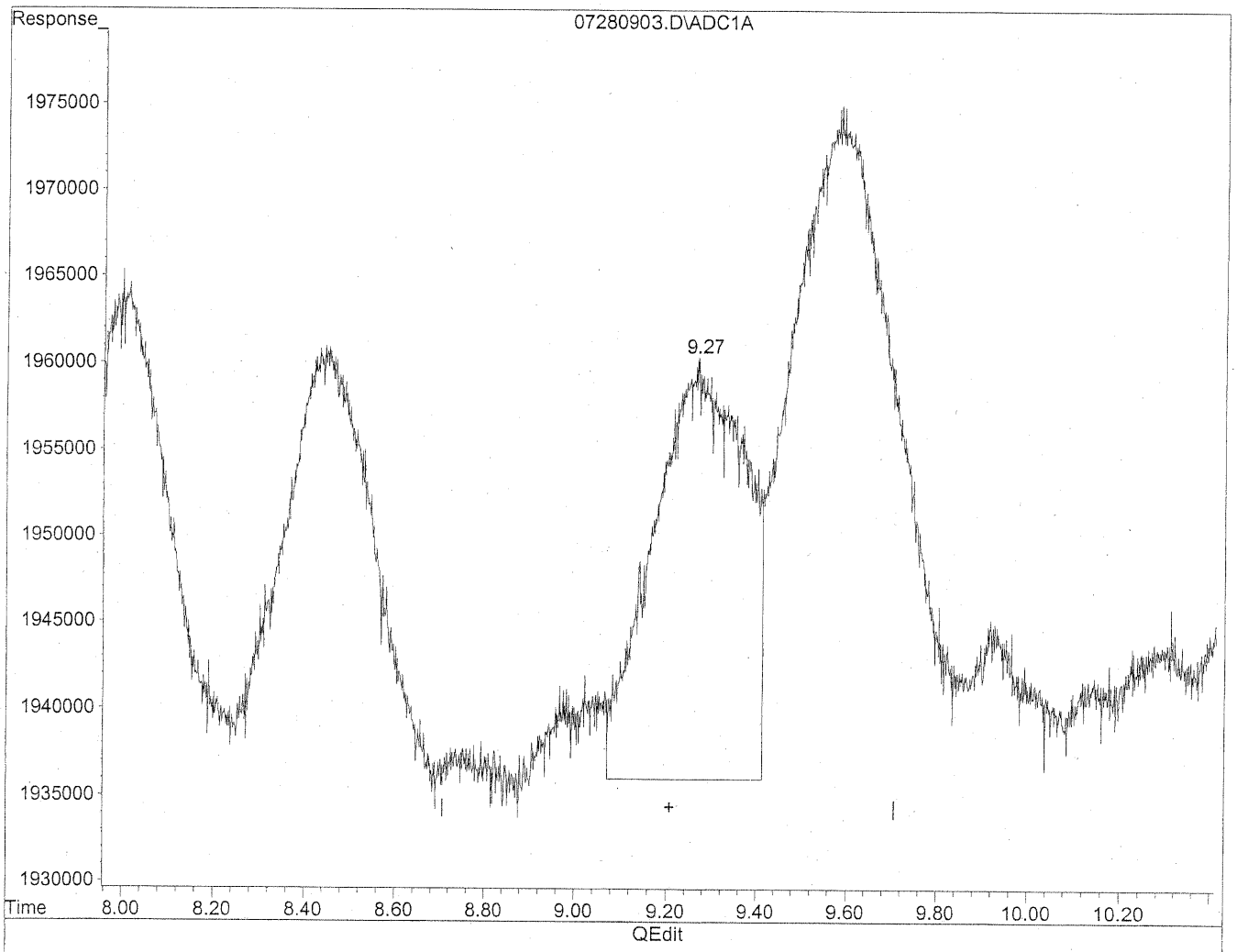
381

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280903.D Vial: 3
Acq On : 28 Jul 2009 9:09 am Operator: HC
Sample : 50ng/ml TO11A Std S21-07270908 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 15:29:52 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.27min 62.877ng/ml m

response 3387183

all factors 1.C

KE 7/29/09

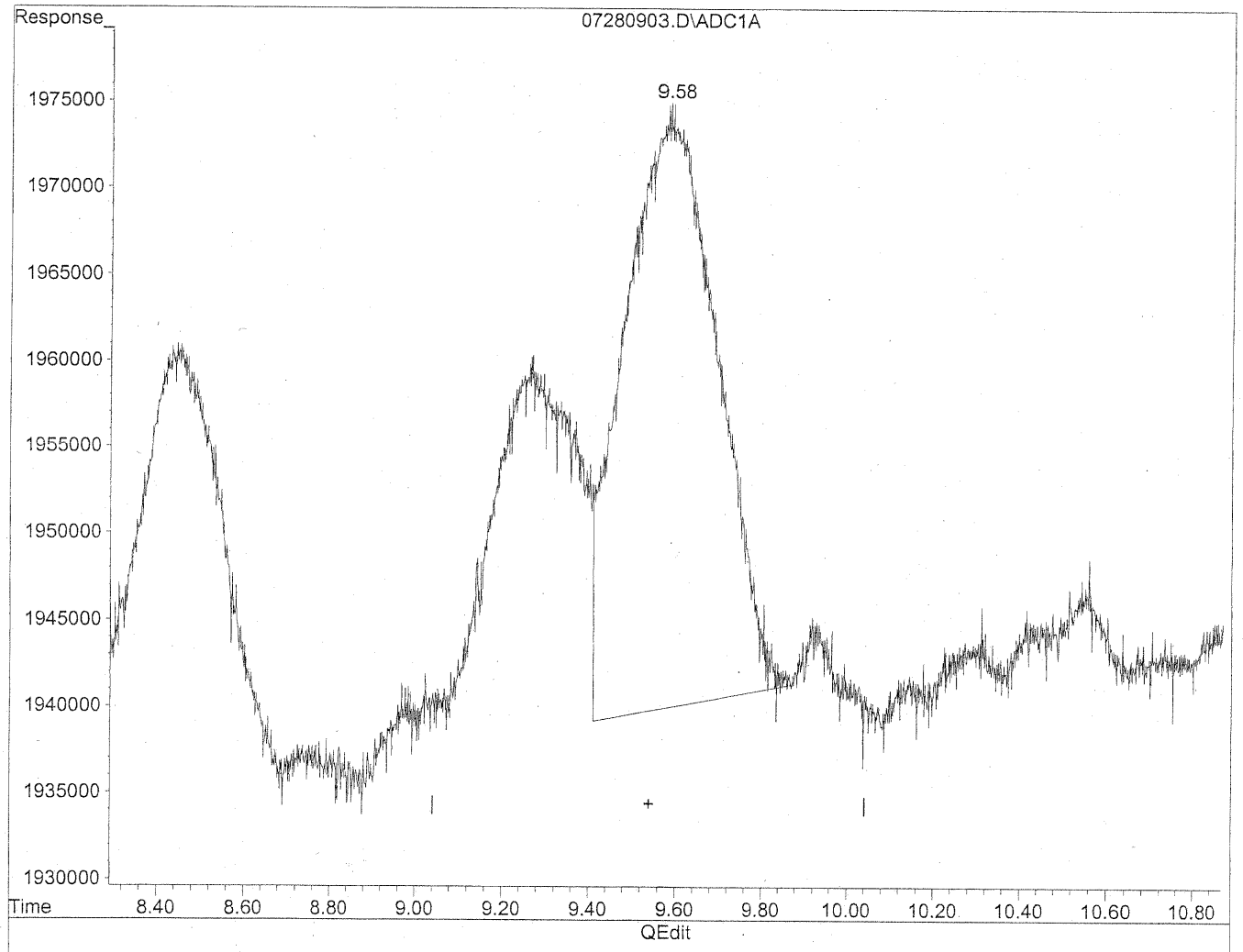
382

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280903.D Vial: 3
Acq On : 28 Jul 2009 9:09 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

9.58min 95.567ng/ml

response 5147699

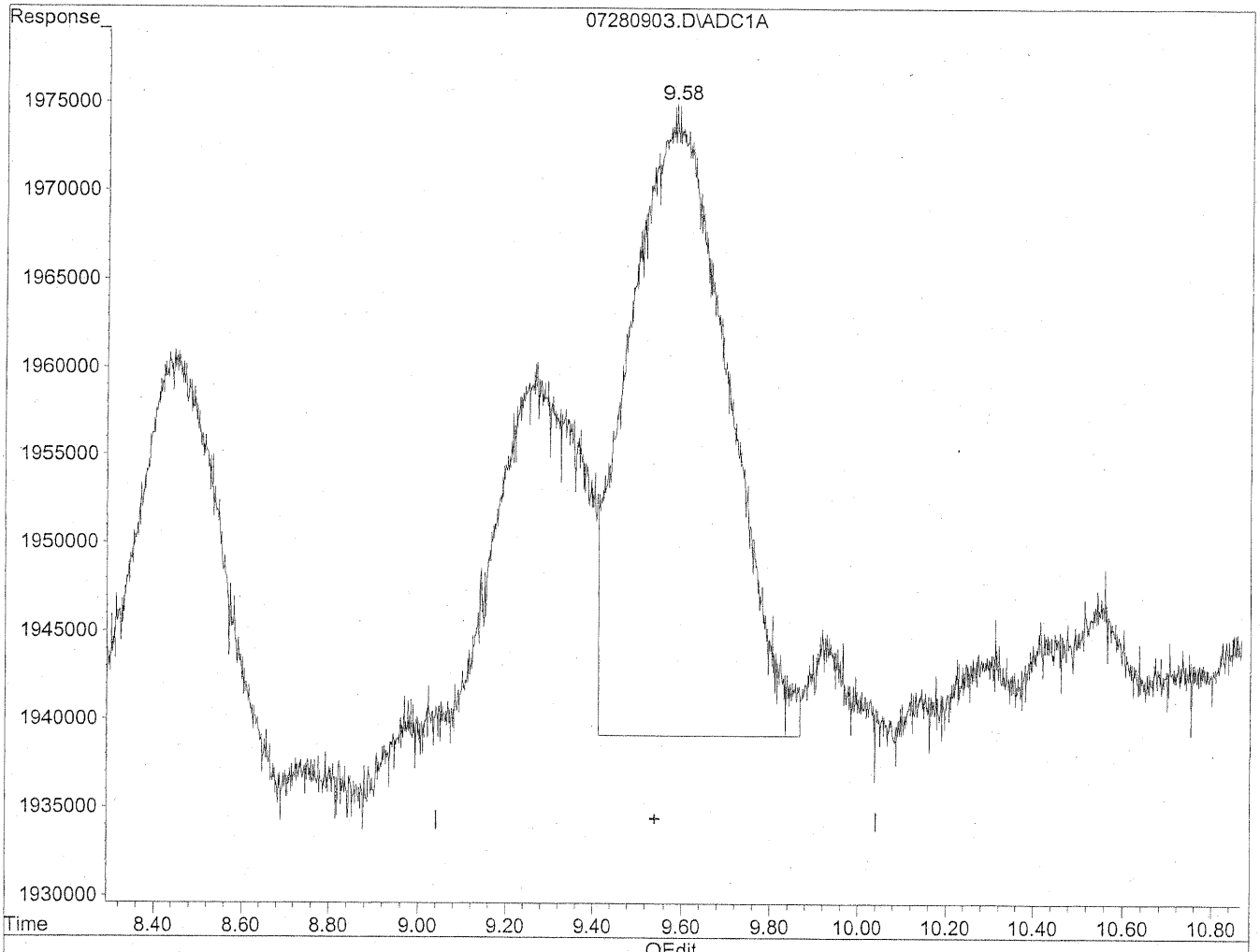
383

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280903.D Vial: 3
Acq On : 28 Jul 2009 9:09 am Operator: HC
Sample : 50ng/ml TO11A Std S21-07270908 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 15:29:52 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

9.58min 101.089ng/ml m

response 5445142

JHC 7/28/09 BC

KR 7/29/09

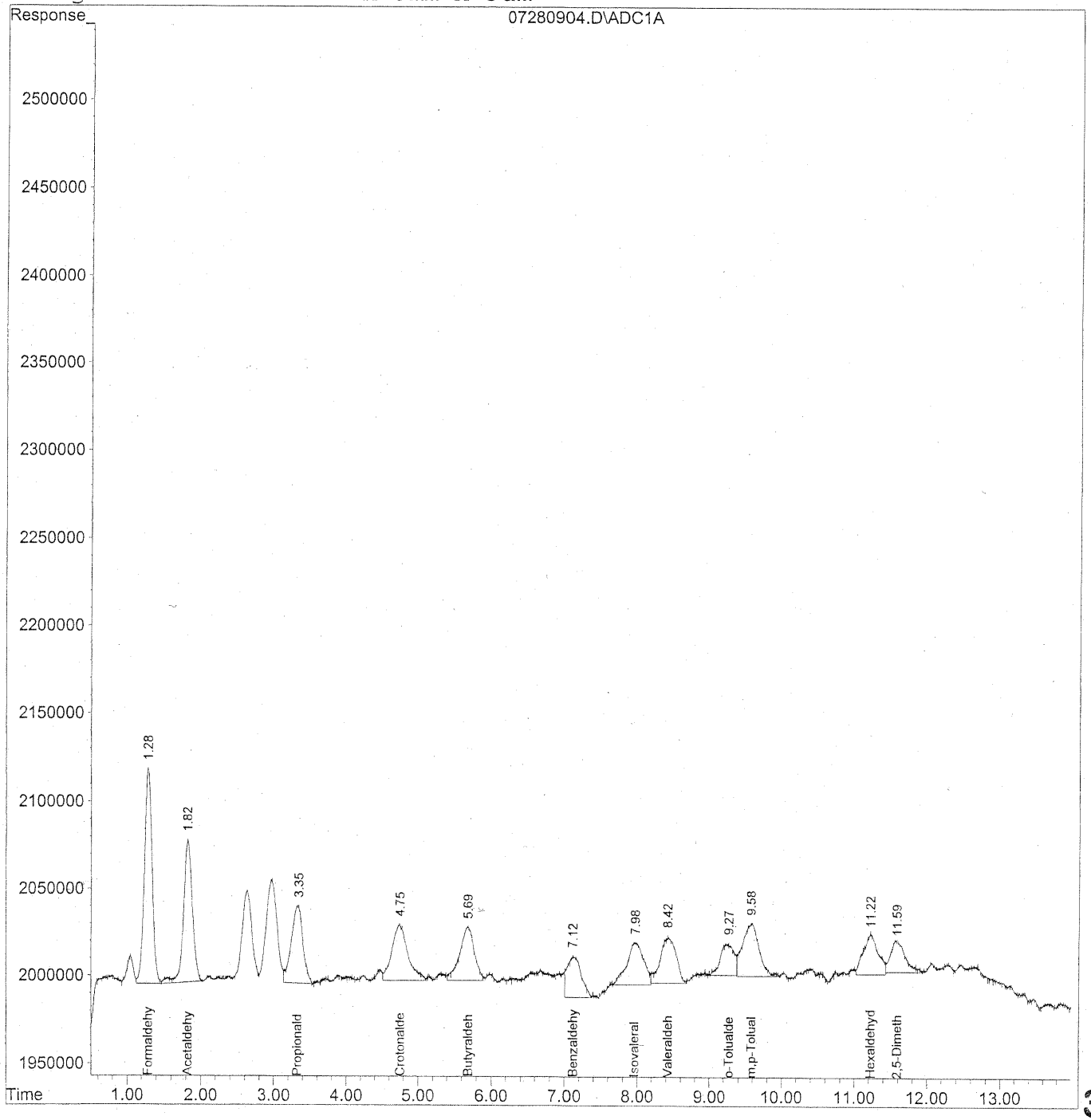
384

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std S21-07270908 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 30 7:38 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std S21-07270908 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 30 7:38 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

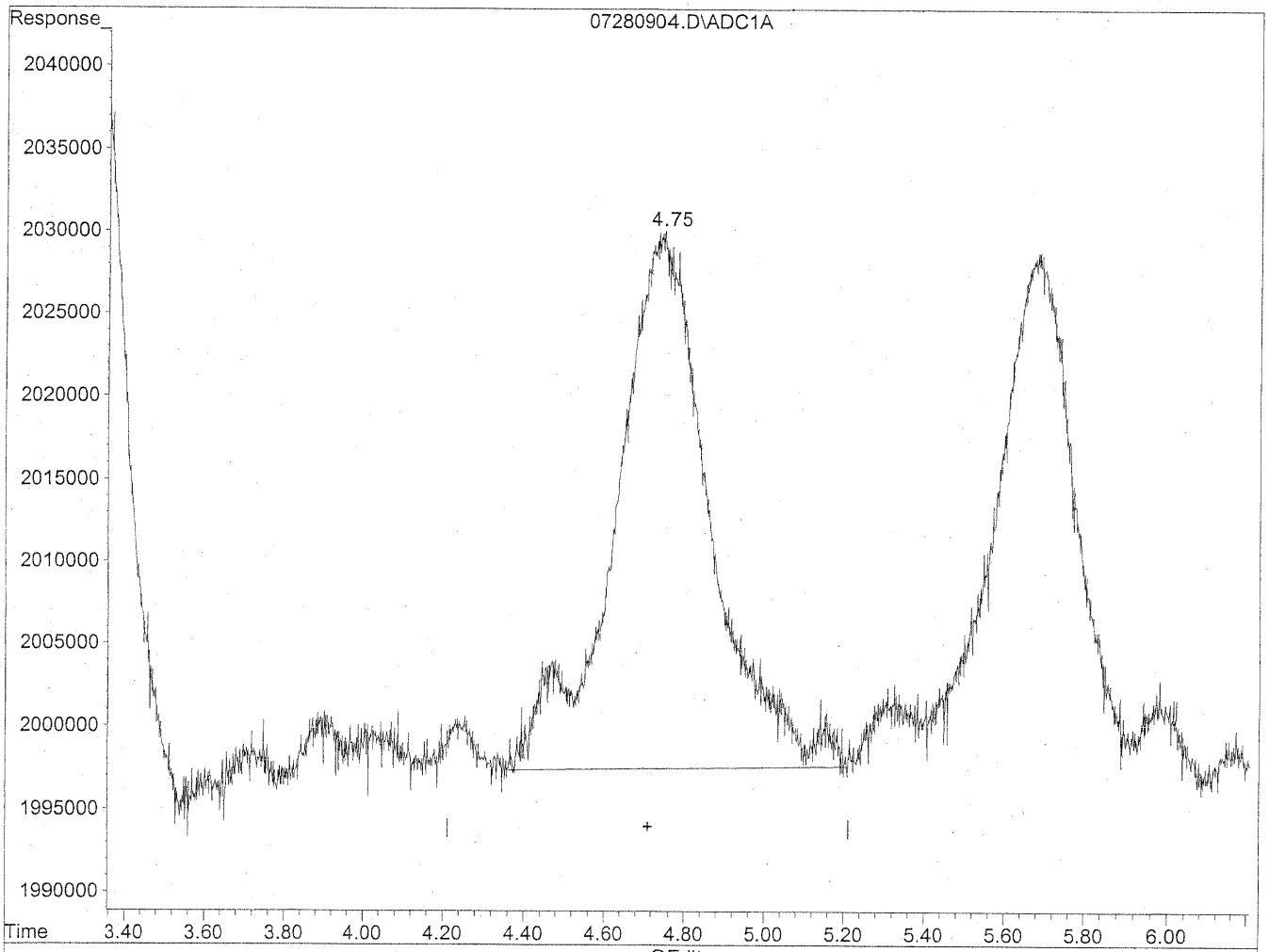
	Compound	R.T.	Response	Conc Units
	Target Compounds			
1)	Formaldehyde	1.29	8859457	50.455 ng/ml
2)	Acetaldehyde	1.82	6975740	51.711 ng/ml
3)	Propionaldehyde	3.33	4973947	48.387 ng/ml
4)	Crotonaldehyde	4.75	4974991	45.000 ng/mlm
5)	Butyraldehyde	5.69	4293221	53.348 ng/mlm
6)	Benzaldehyde	7.12	3079204	48.820 ng/mlm
7)	Isovaleraldehyde	7.96	4002738	45.151 ng/mlm
8)	Valeraldehyde	8.42	4025564	48.445 ng/mlm
9)	o-Tolualdehyde	9.27	2461625	45.695 ng/mlm
10)	m,p-Tolualdehyde	9.58	4897087	90.915 ng/mlm
11)	Hexaldehyde	11.22	3295067	49.079 ng/mlm
12)	2,5-Dimethylbenzaldehyde	11.59	2605446	50.169 ng/mlm

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.74min 48.324ng/ml

response 5342434

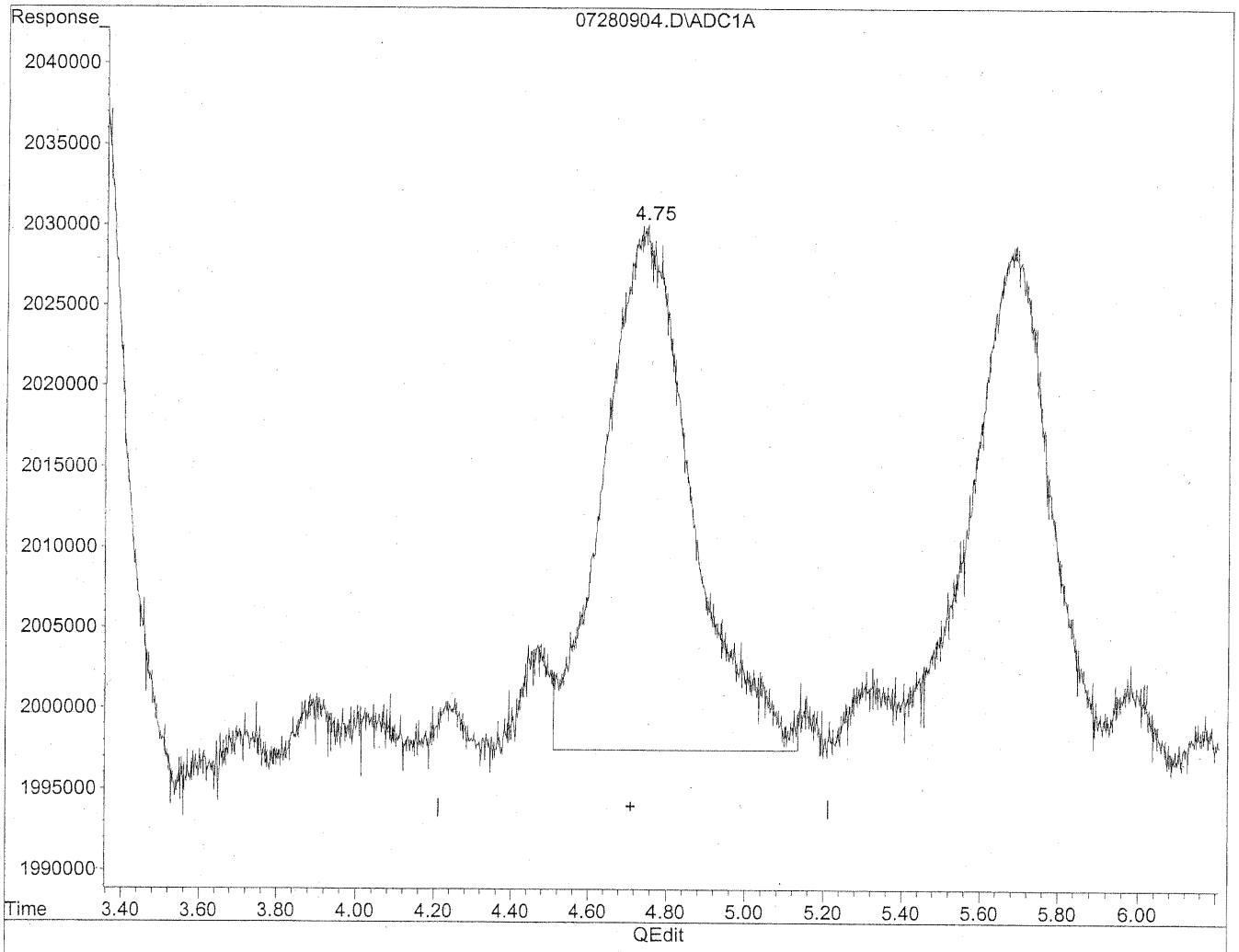
387

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.75min 45.000ng/ml m

response 4974991

HL 7/28/09 CH

KH 7/29/09

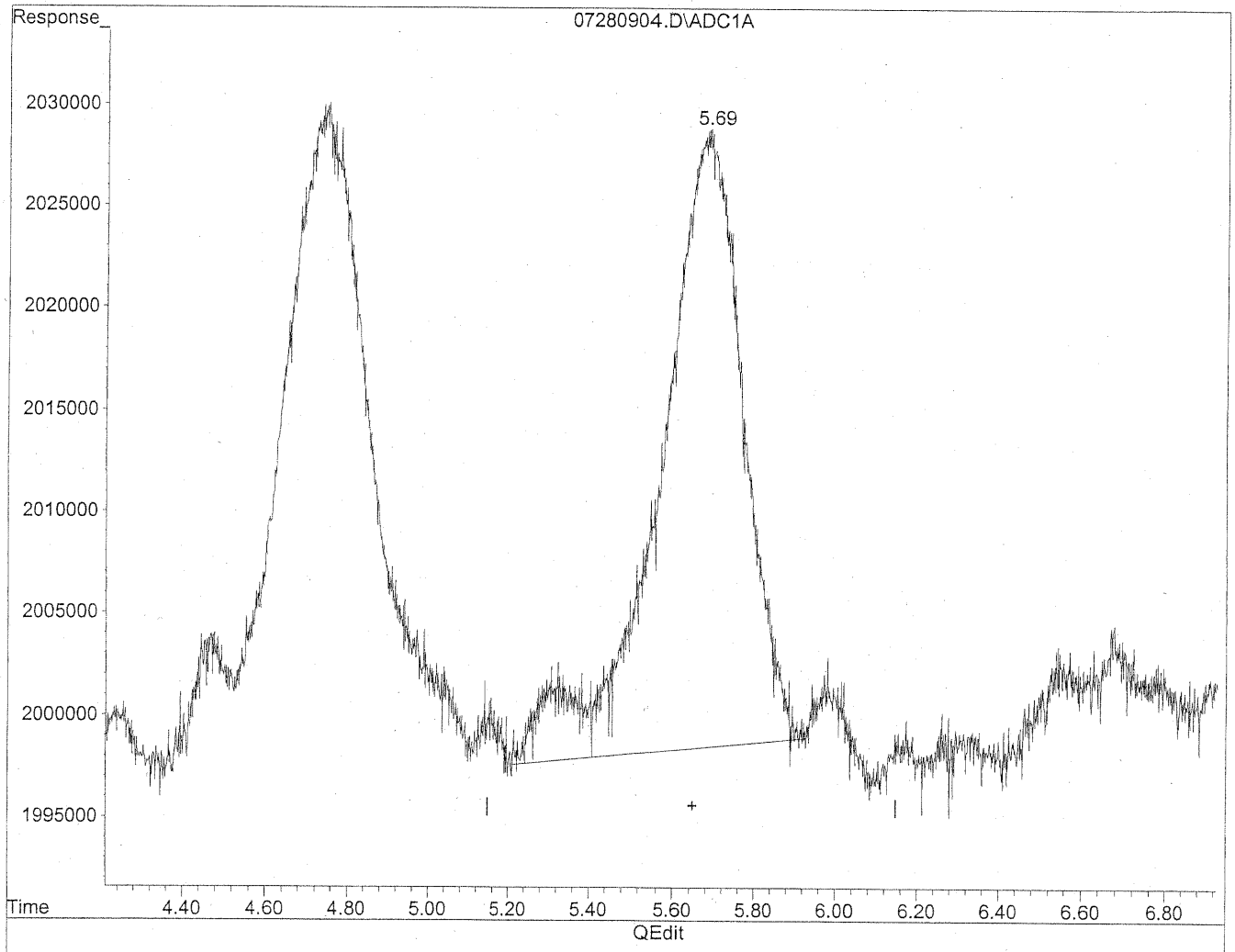
388

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.68min 53.153ng/ml

response 4277470

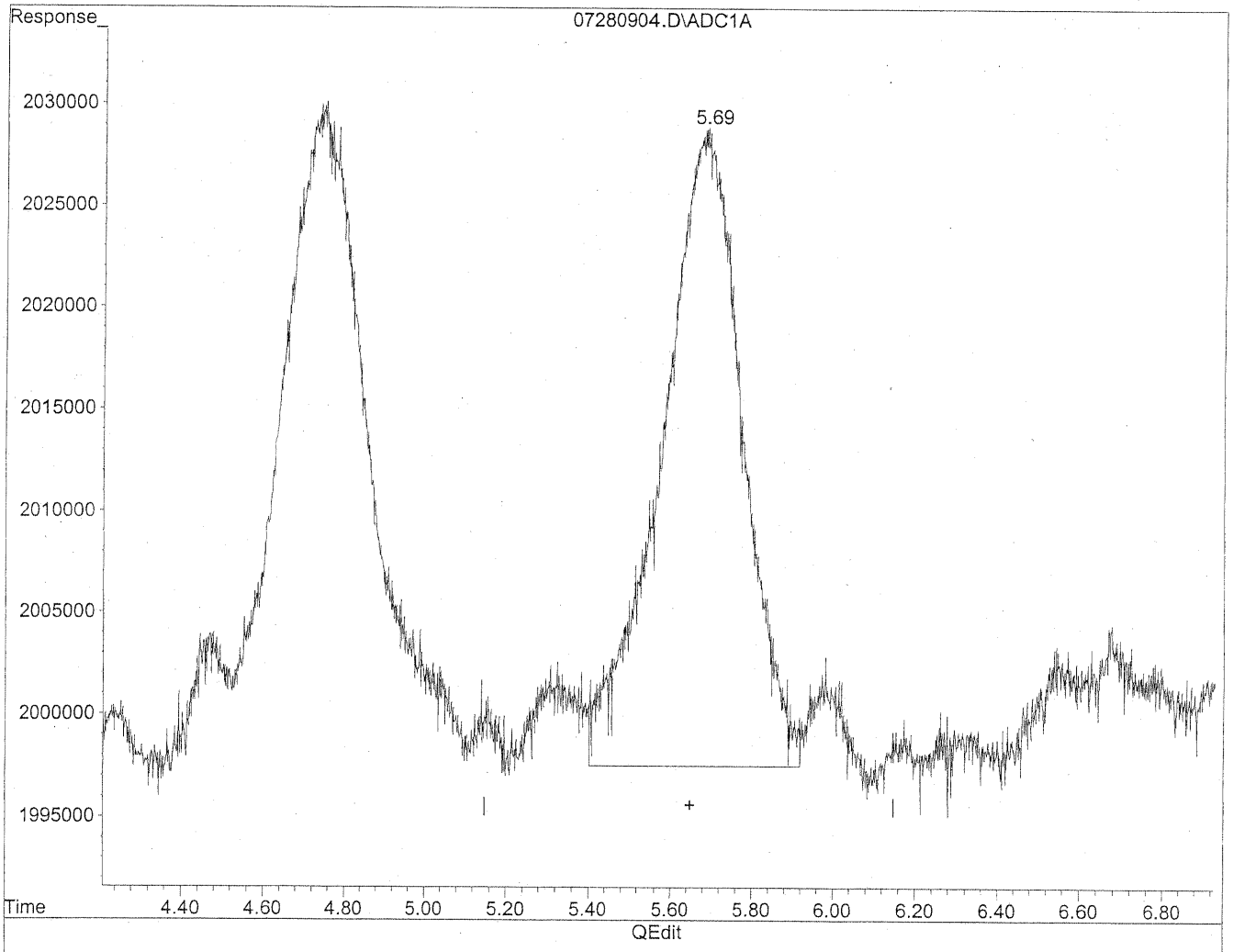
(+) = Expected Retention Time

07280904.D TO110709.M Tue Jul 28 14:23:44 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.69min 53.348ng/ml m

response 4293221

HC 8/28/09 SH

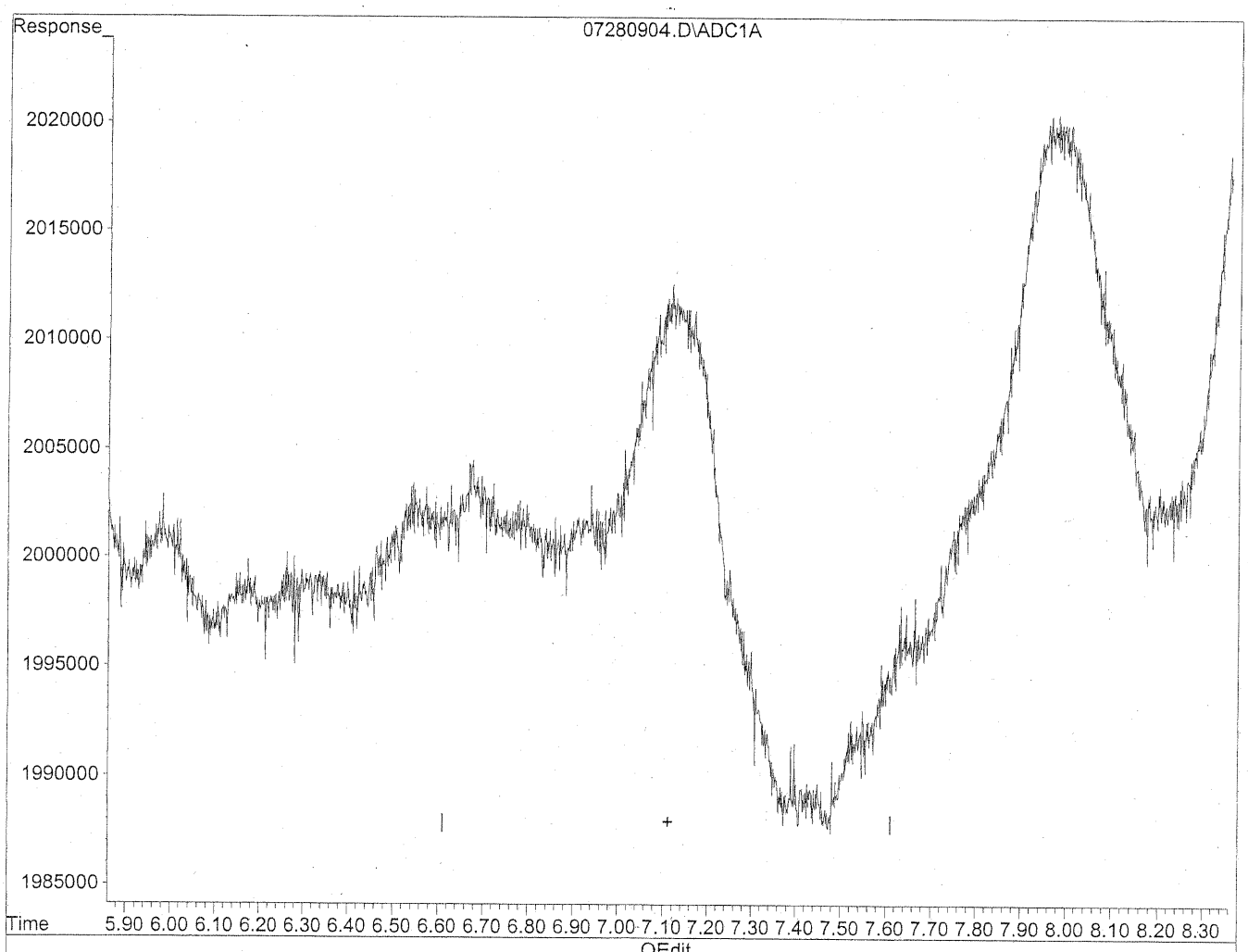
KR 8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

7.11min 0.000ng/ml

response 0

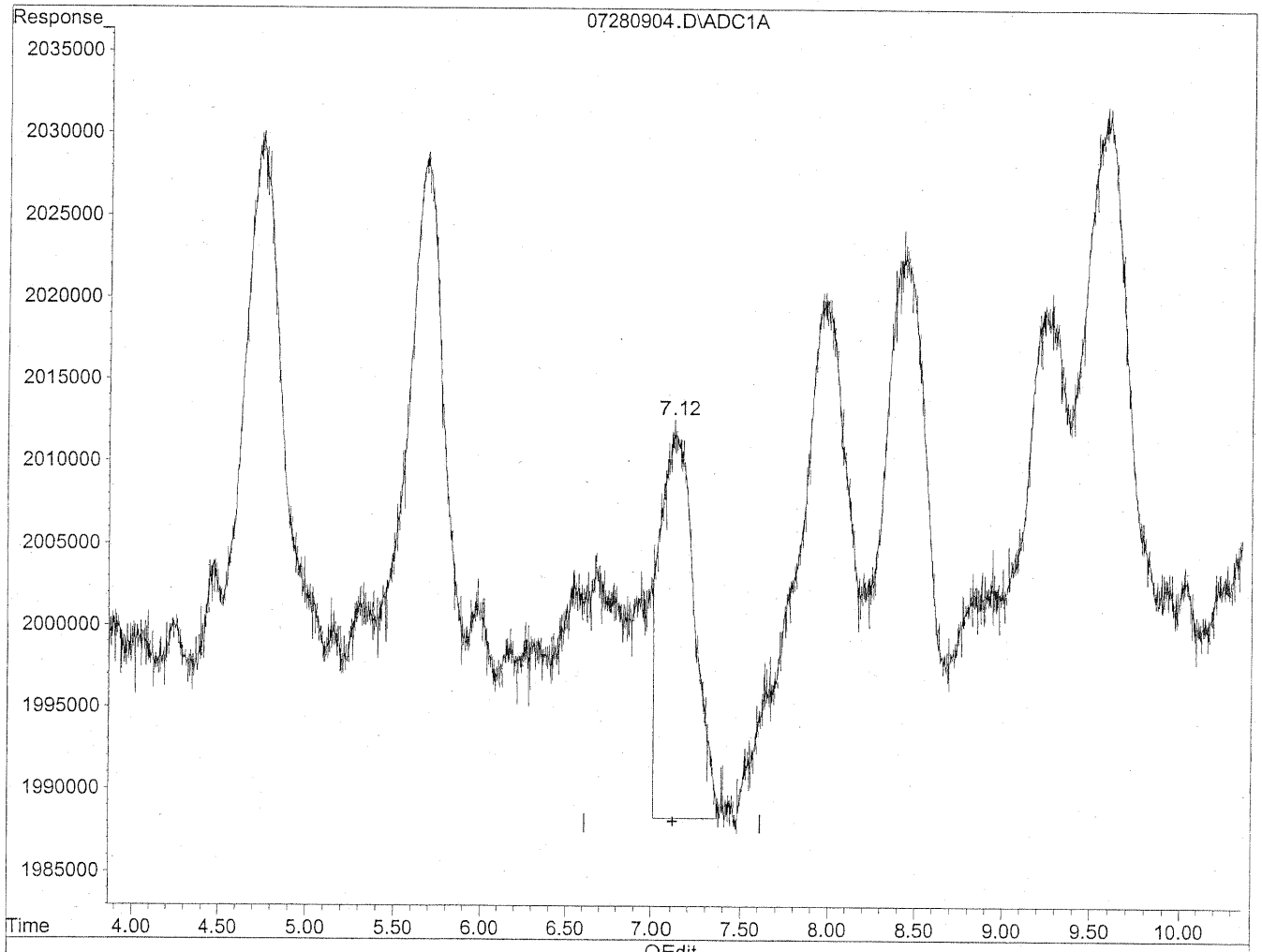
391

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

7.12min 48.820ng/ml m

response 3079204

HC 7/28/09 BW1

KR 7/29/09

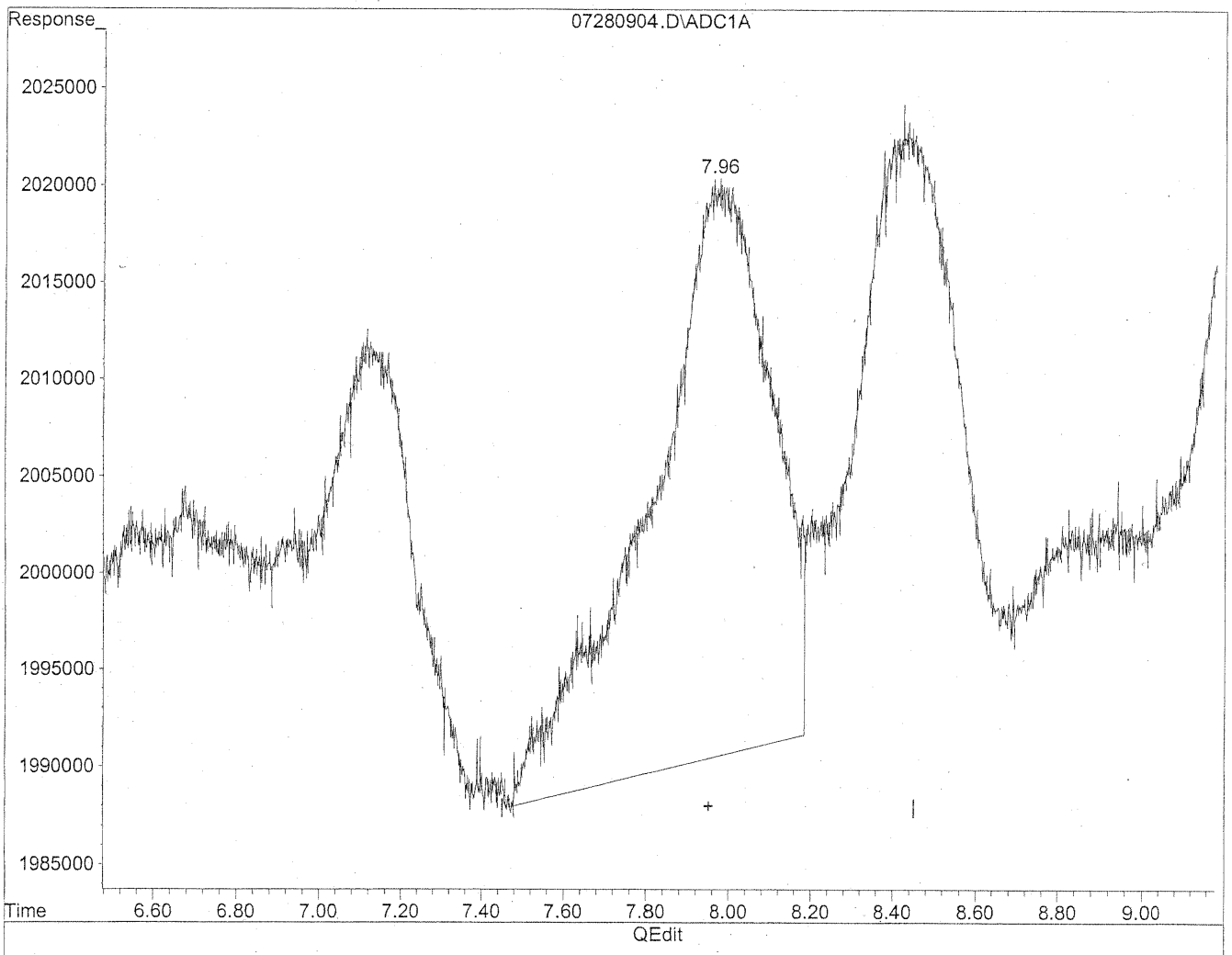
392

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.97min 68.251ng/ml

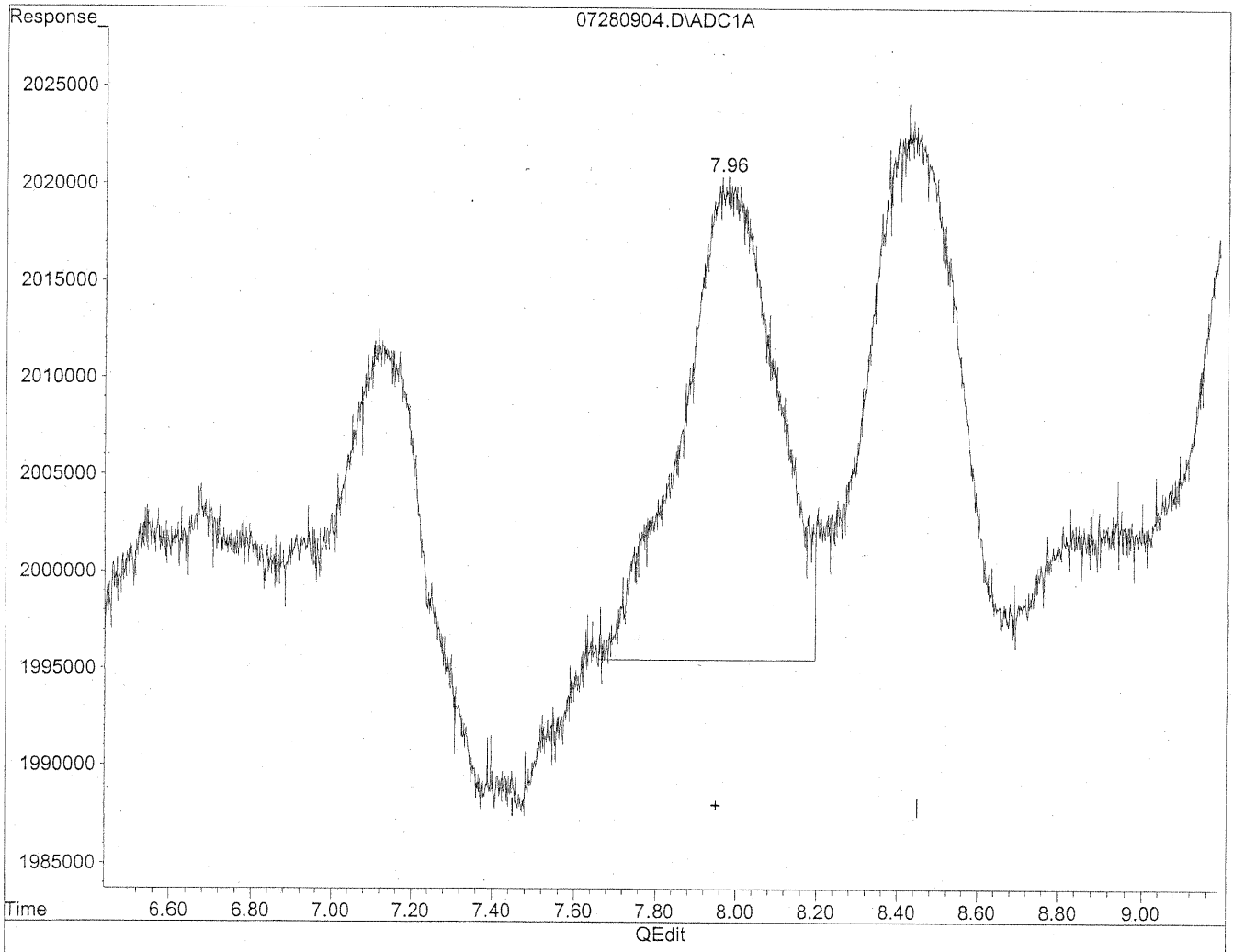
response 6050534

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std S21-07270908 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:25 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Mar 21 12:19:47 2005
Response via : Multiple Level Calibration



QEdit

(7) Isovaleraldehyde

7.96min 45.151ng/ml m

response 4002738

HC 7/29/09
LC

HC 7/30/09

394

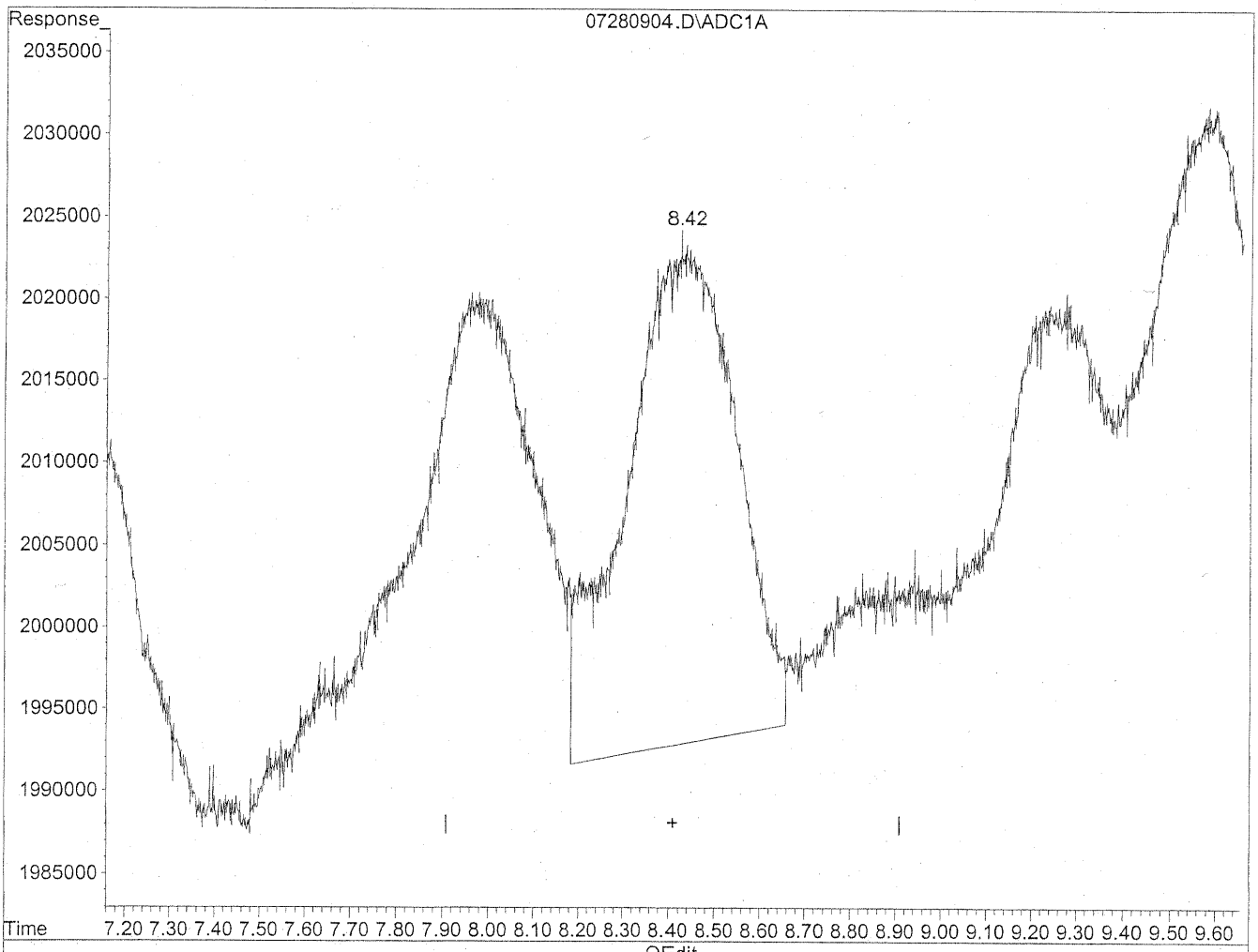
(+) = Expected Retention Time

07280904.D TO110709.M Thu Jul 30 07:38:22 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

8.43min 61.279ng/ml

response 5091976

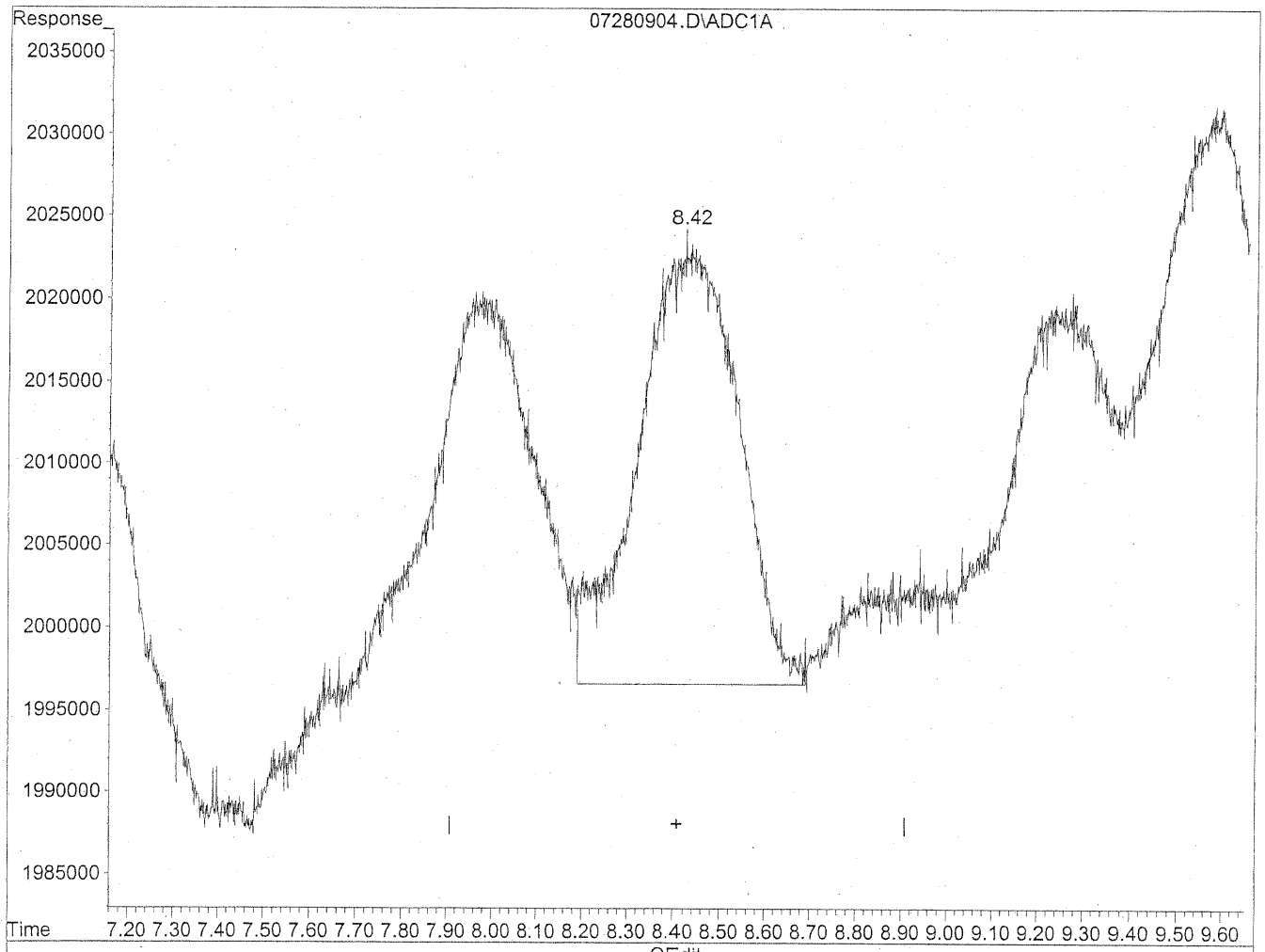
(+) = Expected Retention Time

07280904.D TO110709.M Tue Jul 28 14:24:38 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



QEdit

(8) Valeraldehyde

8.42min 48.445ng/ml m

response 4025564

HC 7/28/09
LC

KR 7/29/09

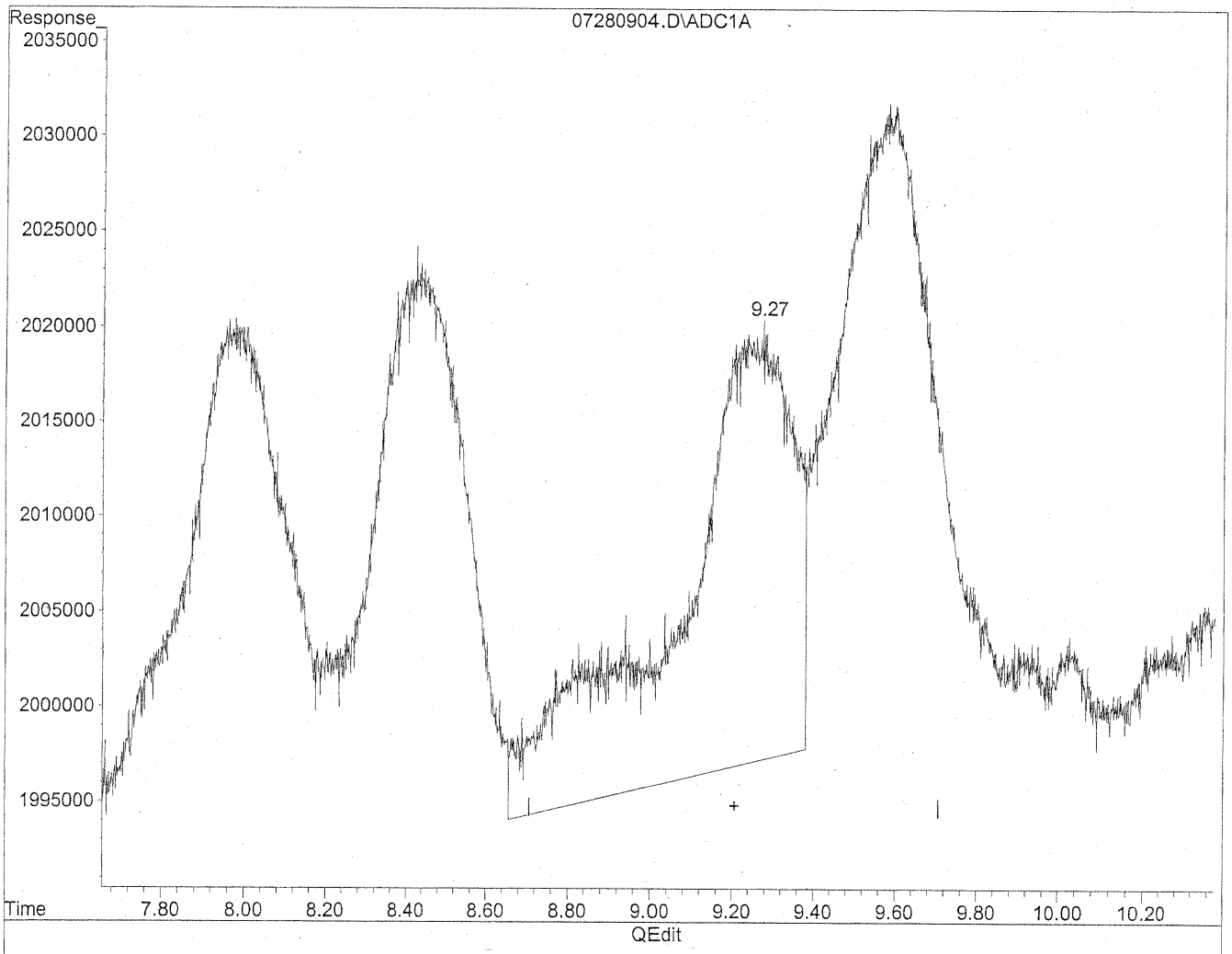
(+) = Expected Retention Time

07280904.D TO110709.M Tue Jul 28 14:24:48 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.24min 84.965ng/ml

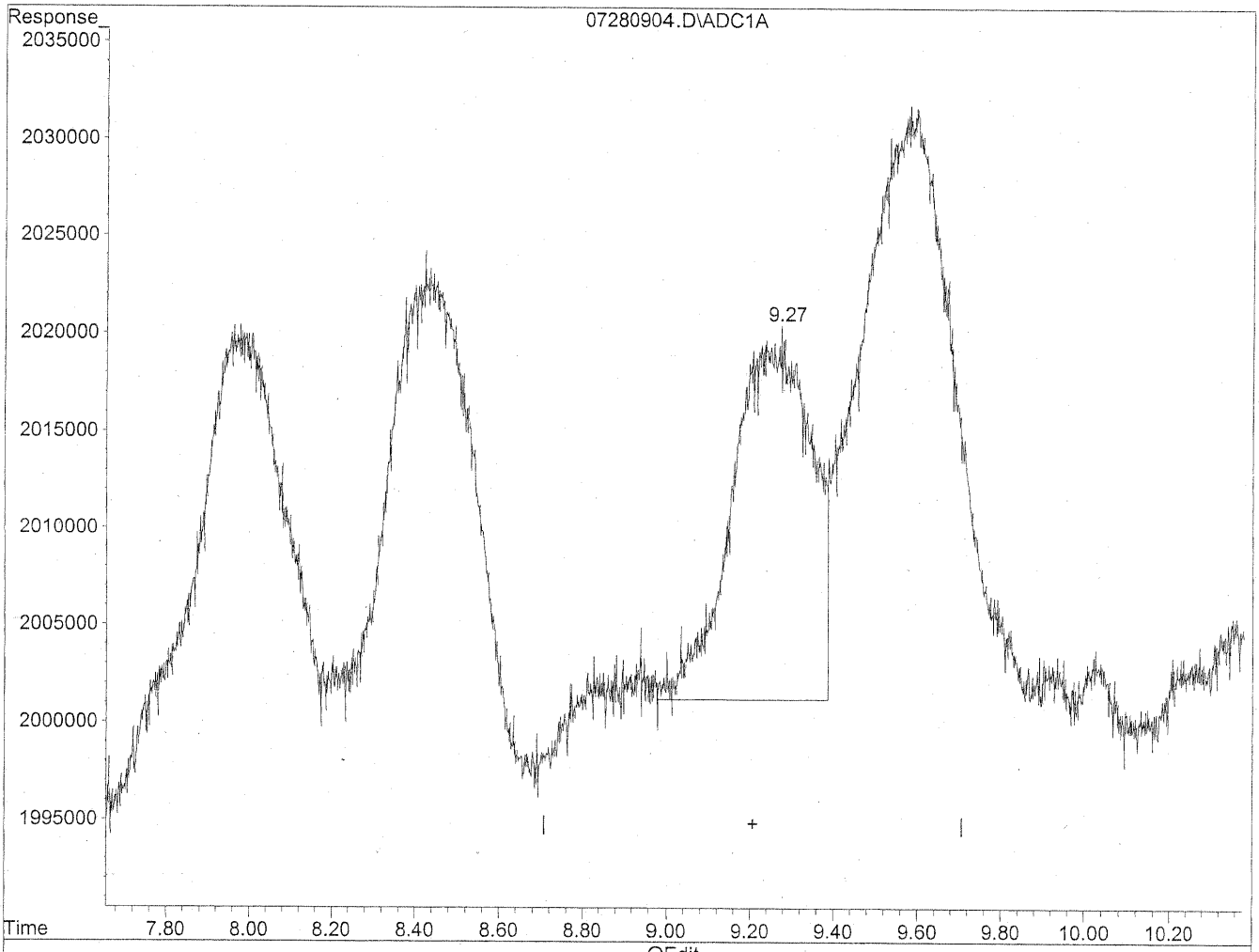
response 4577075

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.27min 45.695ng/ml m

response 2461625

HC 7/28/09 LC

KR 7/29/09

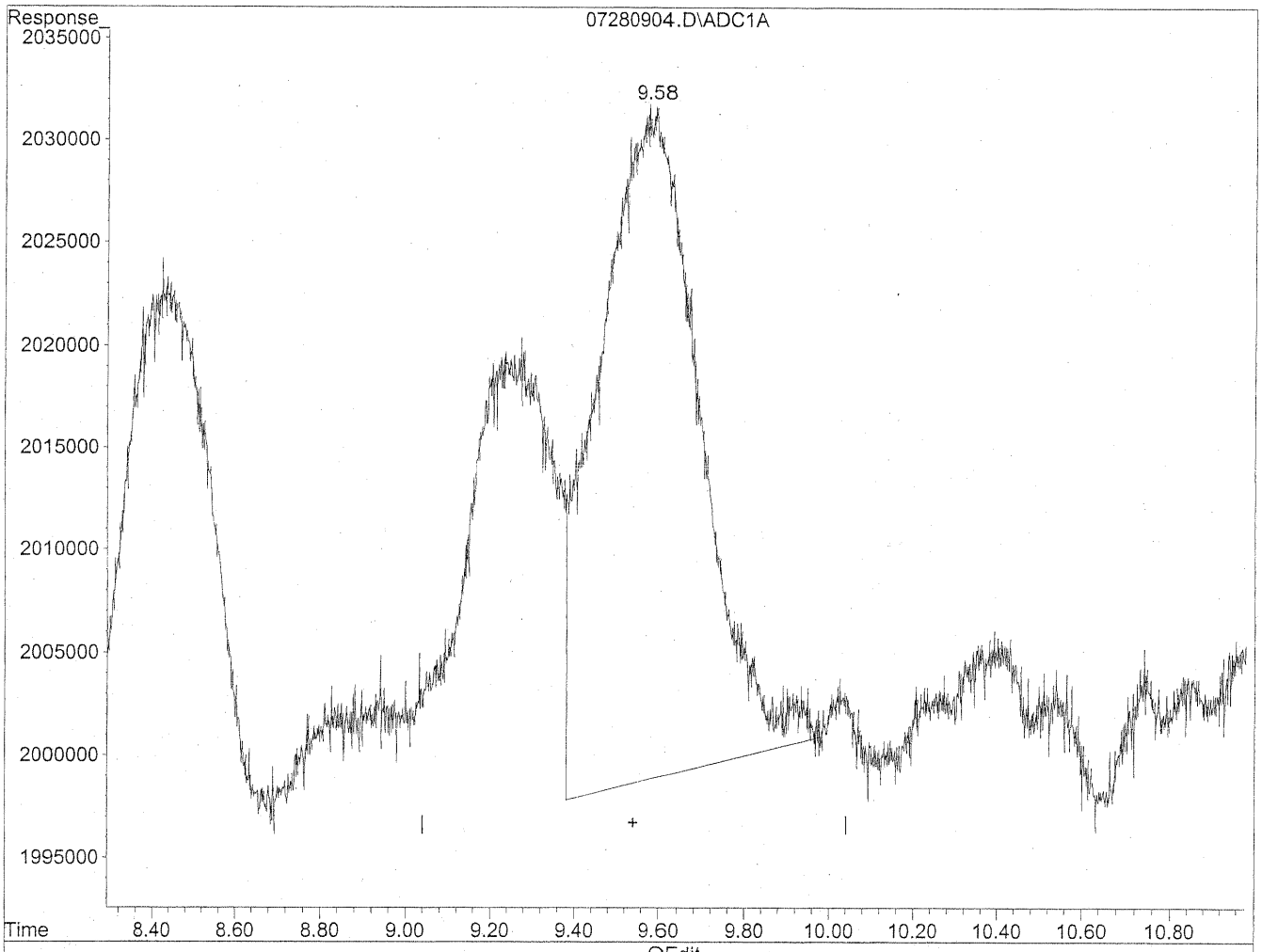
398

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

9.59min 100.987ng/ml

response 5439618

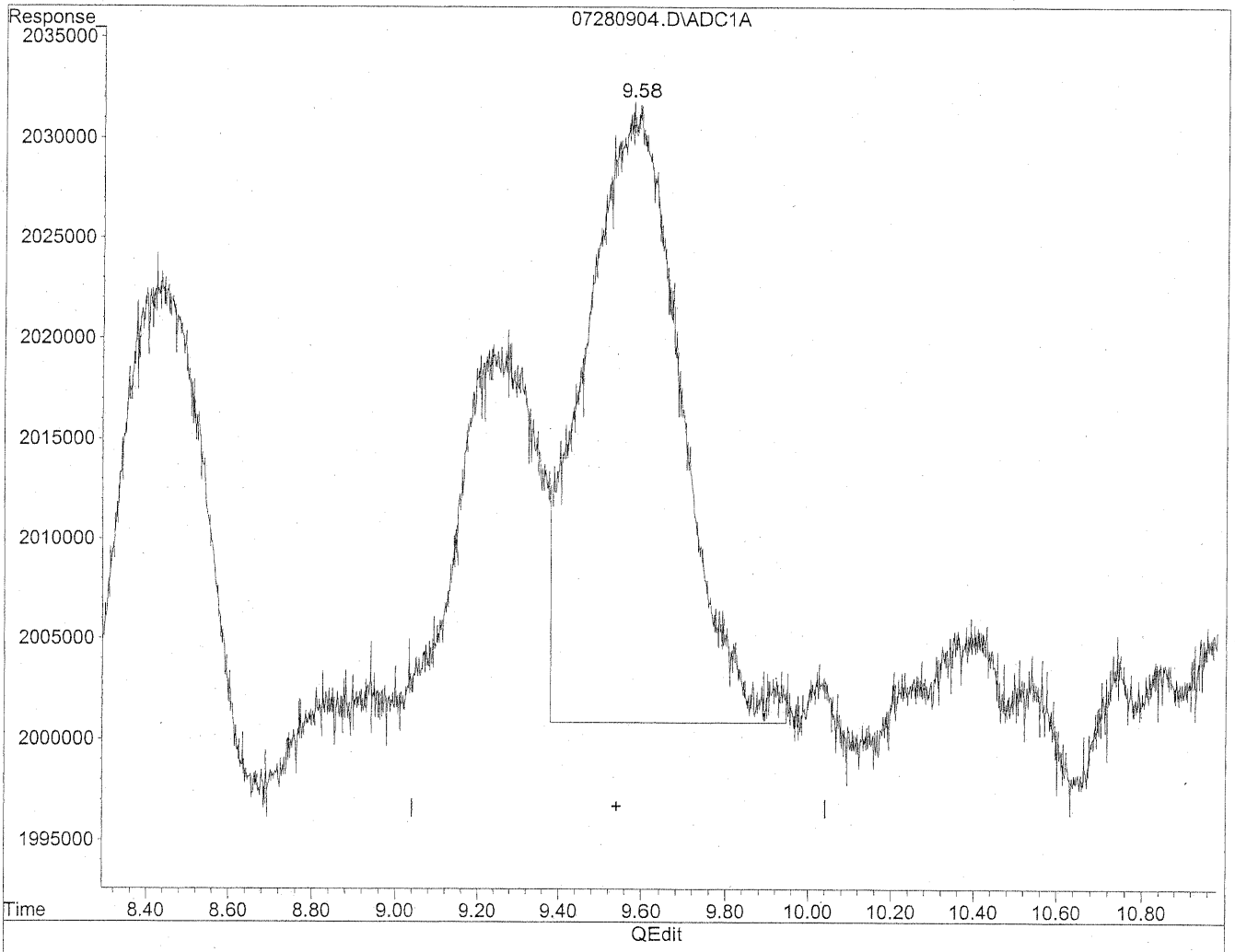
399

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

9.58min 90.915ng/ml m

response 4897087

HC 7/28/09

KR 7/29/09

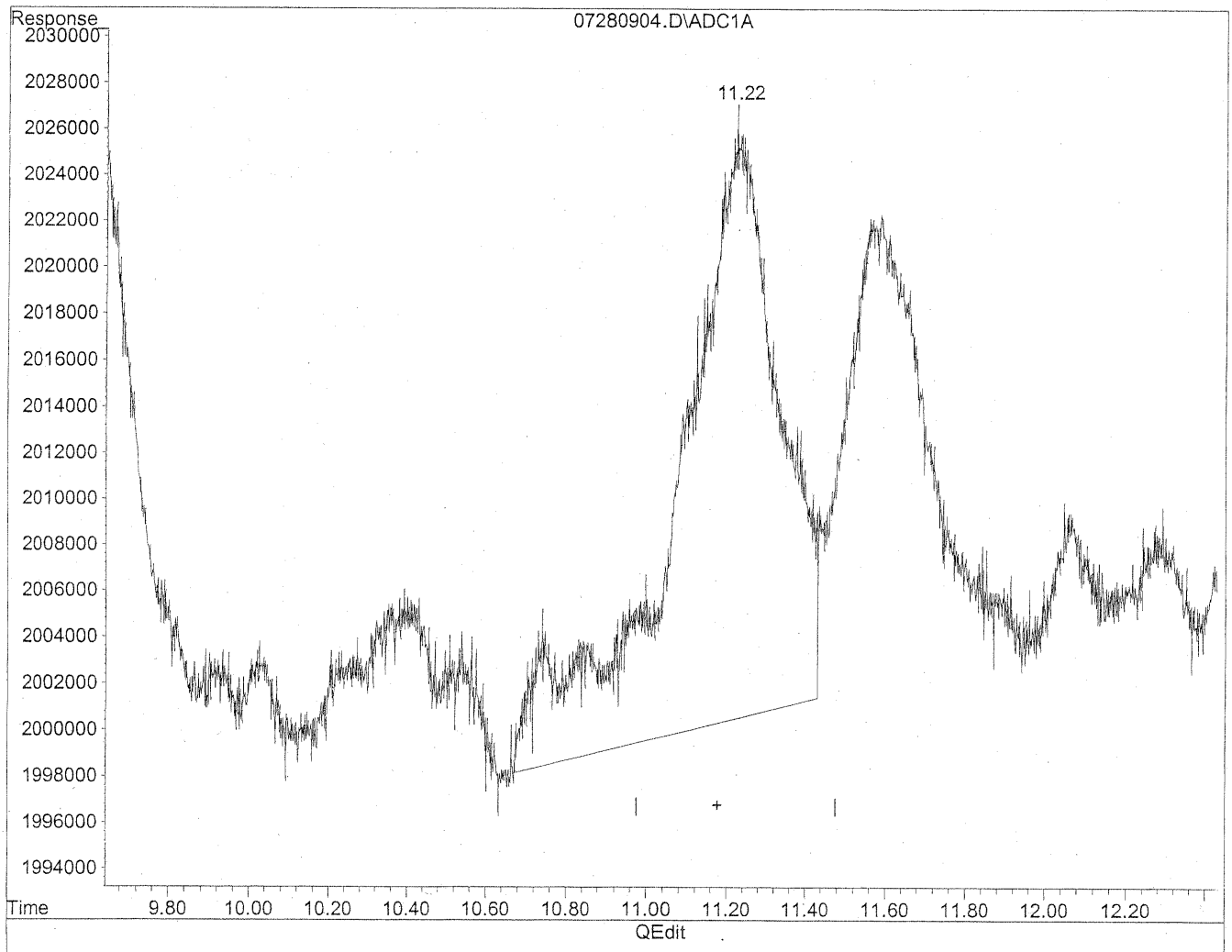
400

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

11.23min 66.912ng/ml

response 4492347

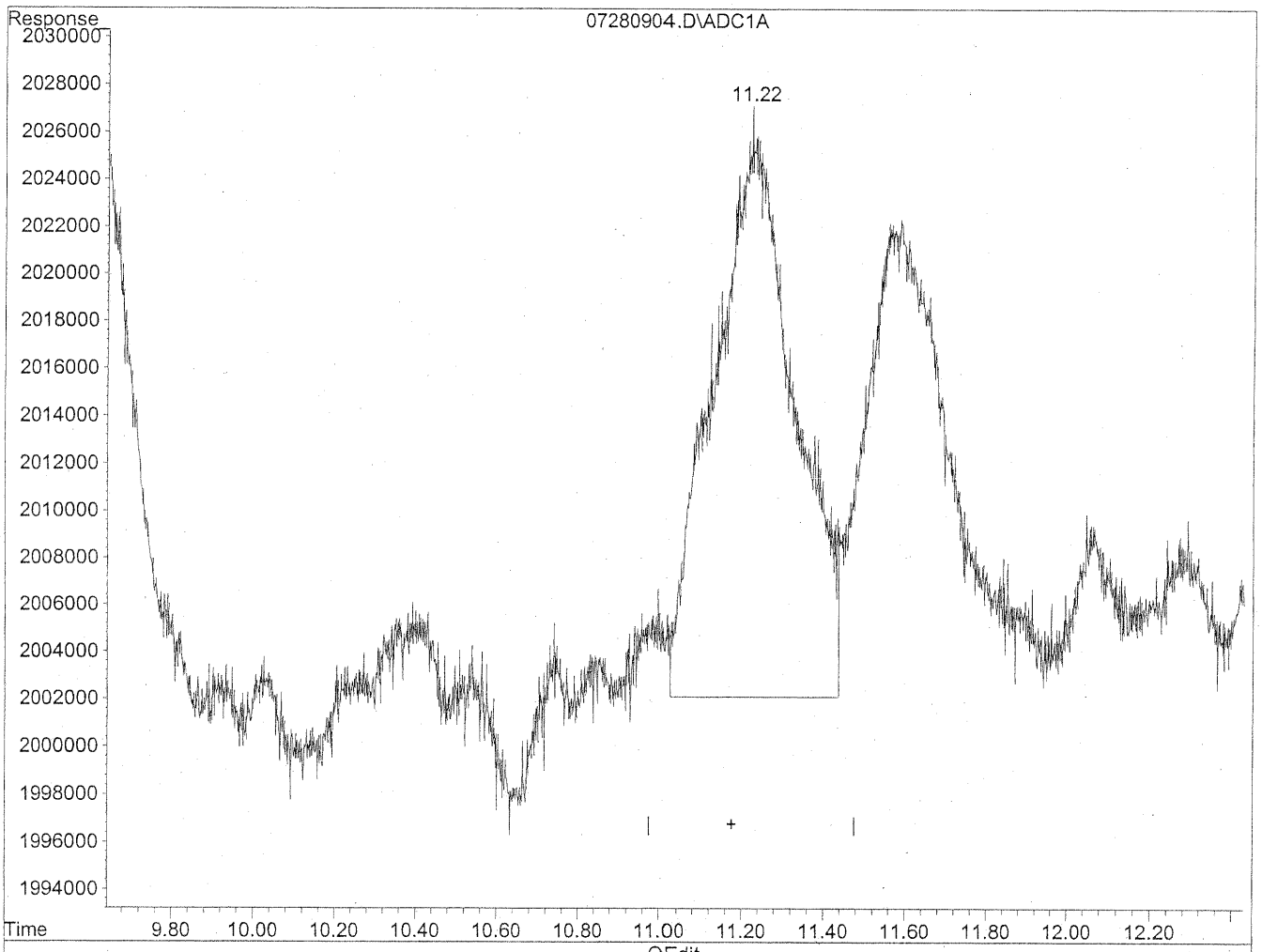
401

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

11.22min 49.079ng/ml m

response 3295067

HC 7/28/09 SH

KL 7/29/09

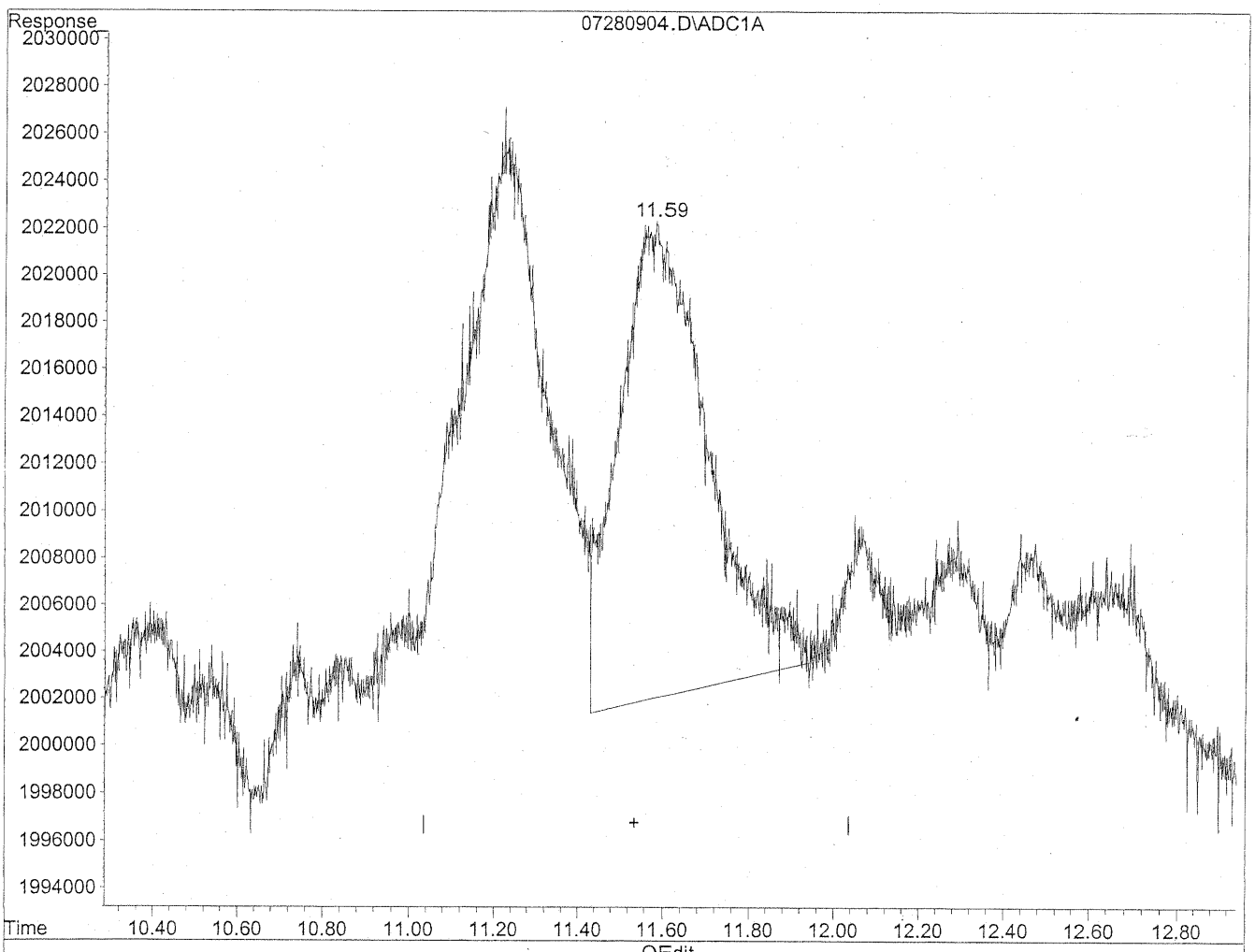
(+) = Expected Retention Time

07280904.D TO110709.M Tue Jul 28 14:25:22 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

11.58min 55.789ng/ml

response 2897339

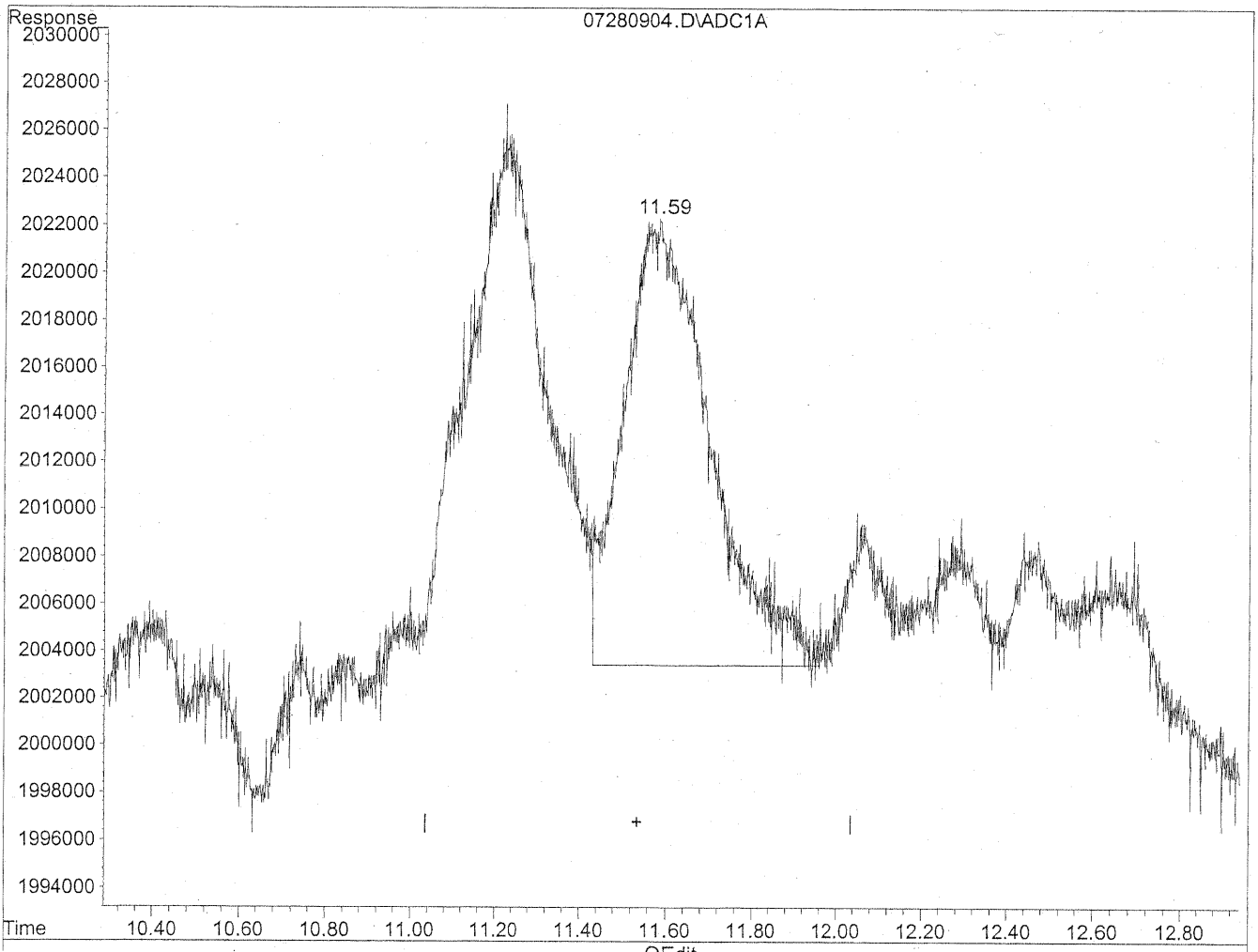
403

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

11.59min 50.169ng/ml m

response 2605446

tc 7/28/09 LC

7/29/09

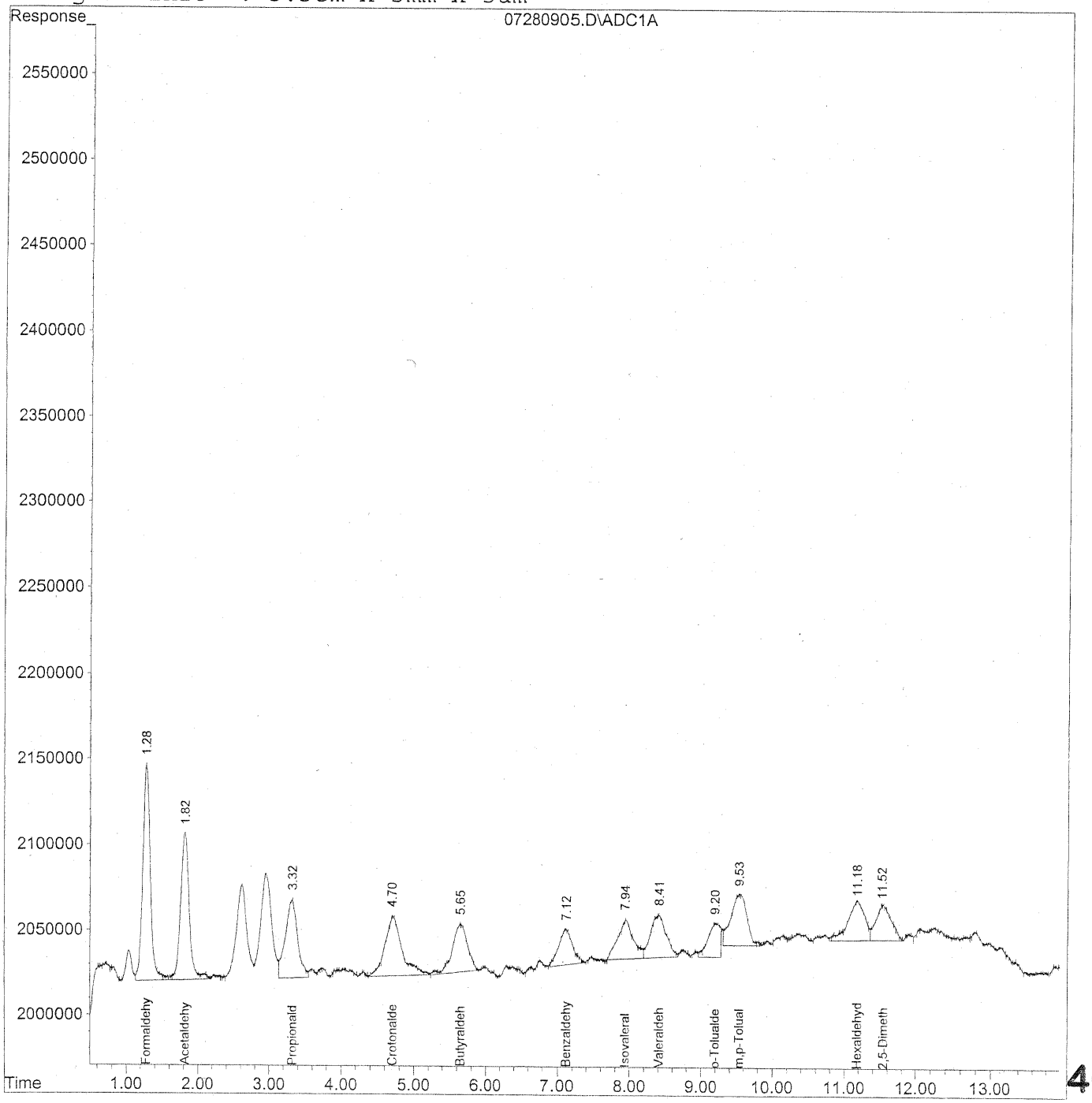
404

(+) = Expected Retention Time
07280904.D TO110709.M Tue Jul 28 14:25:43 2009

Data File : J:\LC01\DATA\TO11\2009_07\28\07280905.D Vial: 5
Acq On : 28 Jul 2009 9:39 am Operator: HC
Sample : 50ng/ml TO11A Std S21-07270908 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:29 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



405

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280905.D Vial: 5
Acq On : 28 Jul 2009 9:39 am Operator: HC
Sample : 50ng/ml TO11A Std S21-07270908 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:29 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.28	9305088	52.992	ng/ml
2)	Acetaldehyde	1.81	7389770	54.780	ng/ml
3)	Propionaldehyde	3.31	5442713	52.947	ng/ml
4)	Crotonaldehyde	4.71	5754474	52.051	ng/ml
5)	Butyraldehyde	5.65	4119144	51.185	ng/ml
6)	Benzaldehyde	7.11	2732056	43.316	ng/ml
7)	Isovaleraldehyde	7.95	3500271	39.483	ng/ml
8)	Valeraldehyde	8.41	3855749	46.402	ng/ml
9)	o-Tolualdehyde	9.20	2416389	44.856	ng/mlm
10)	m,p-Tolualdehyde	9.53	4801019	89.131	ng/mlm
11)	Hexaldehyde	11.18	3739368	55.696	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.54	3118537	60.048	ng/ml

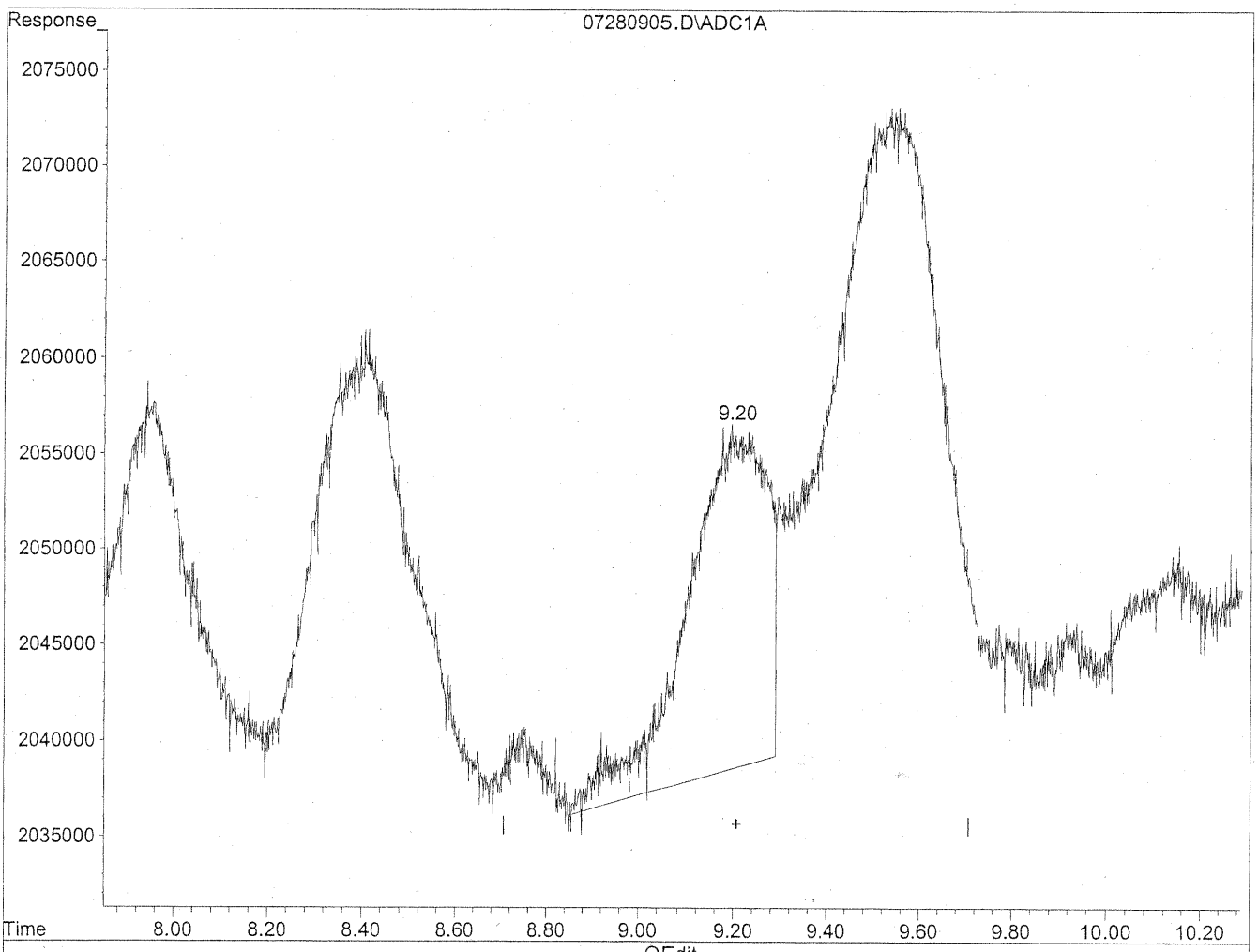
406

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280905.D Vial: 5
Acq On : 28 Jul 2009 9:39 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:25 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.21min 38.587ng/ml

response 2078690

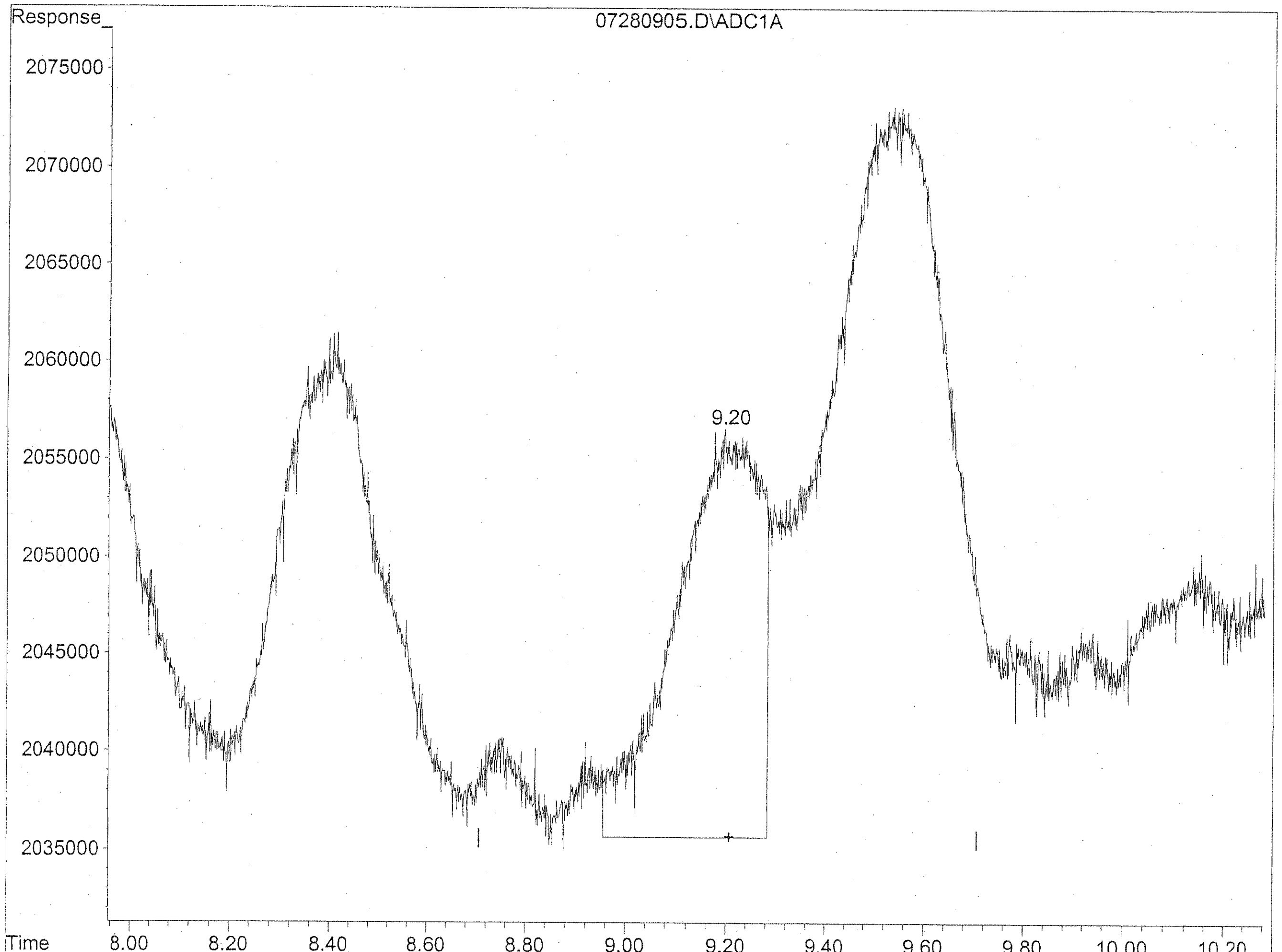
(+) = Expected Retention Time

07280905.D TO110709.M Tue Jul 28 14:26:10 2009

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2009_07\28\07280905.D        Vial: 5
Acq On    : 28 Jul 2009   9:39 am                       Operator: HC
Sample    : 50ng/ml TO11A Std S21-07270908              Inst    : LC 01
Misc      :                                             Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Jul 28 14:29 19109  Quant Results File: TO110709.RES

Method          : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title           : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update     : Tue Jul 28 15:29:52 2009
Response via    : Multiple Level Calibration
```



QEdit

(9) o-Tolualdehyde

9.20min 44.856ng/ml m

response 2416389

HC 7/29/09 IC

KL 7/29/09

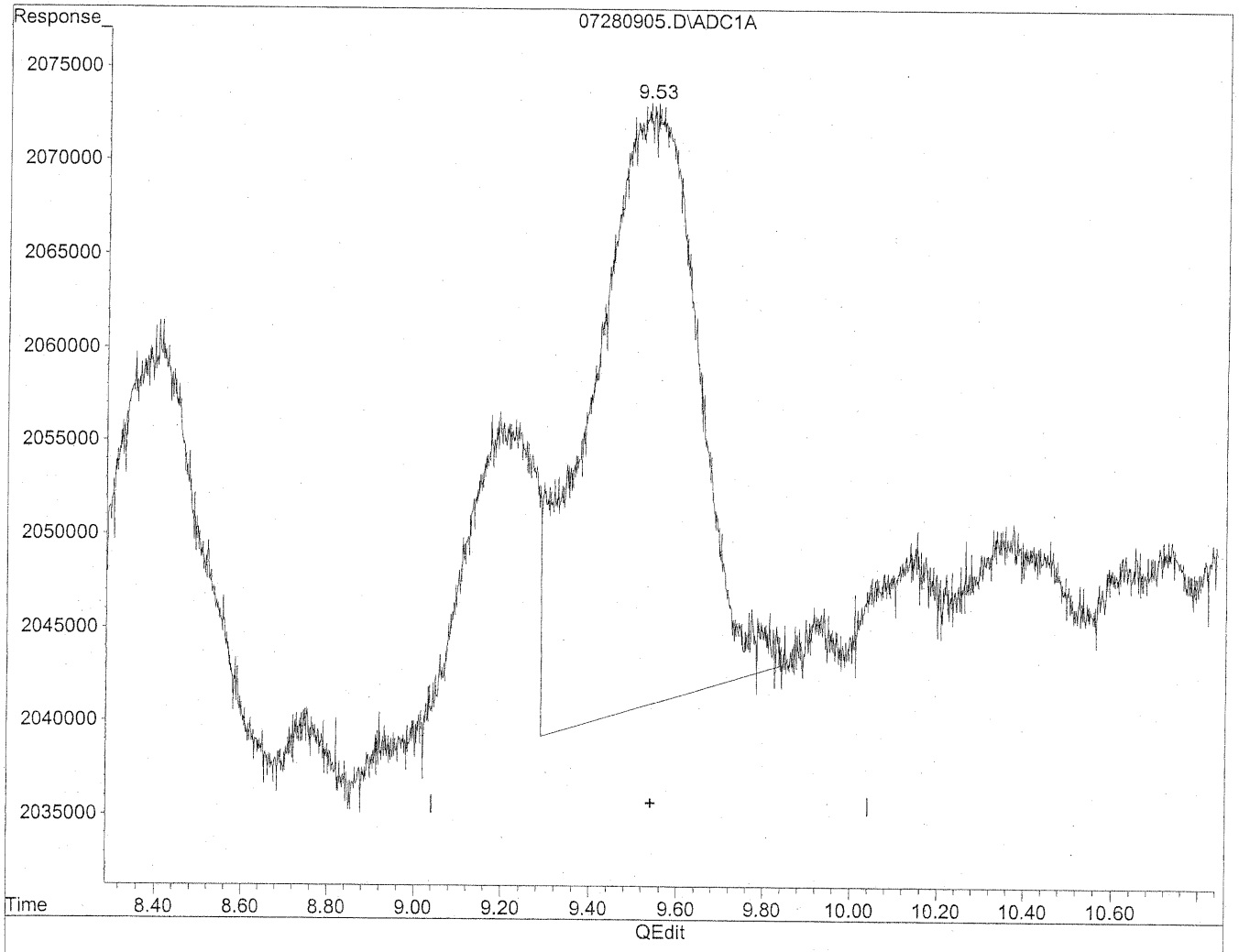
408

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280905.D Vial: 5
Acq On : 28 Jul 2009 9:39 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:25 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

9.54min 100.090ng/ml

response 5391328

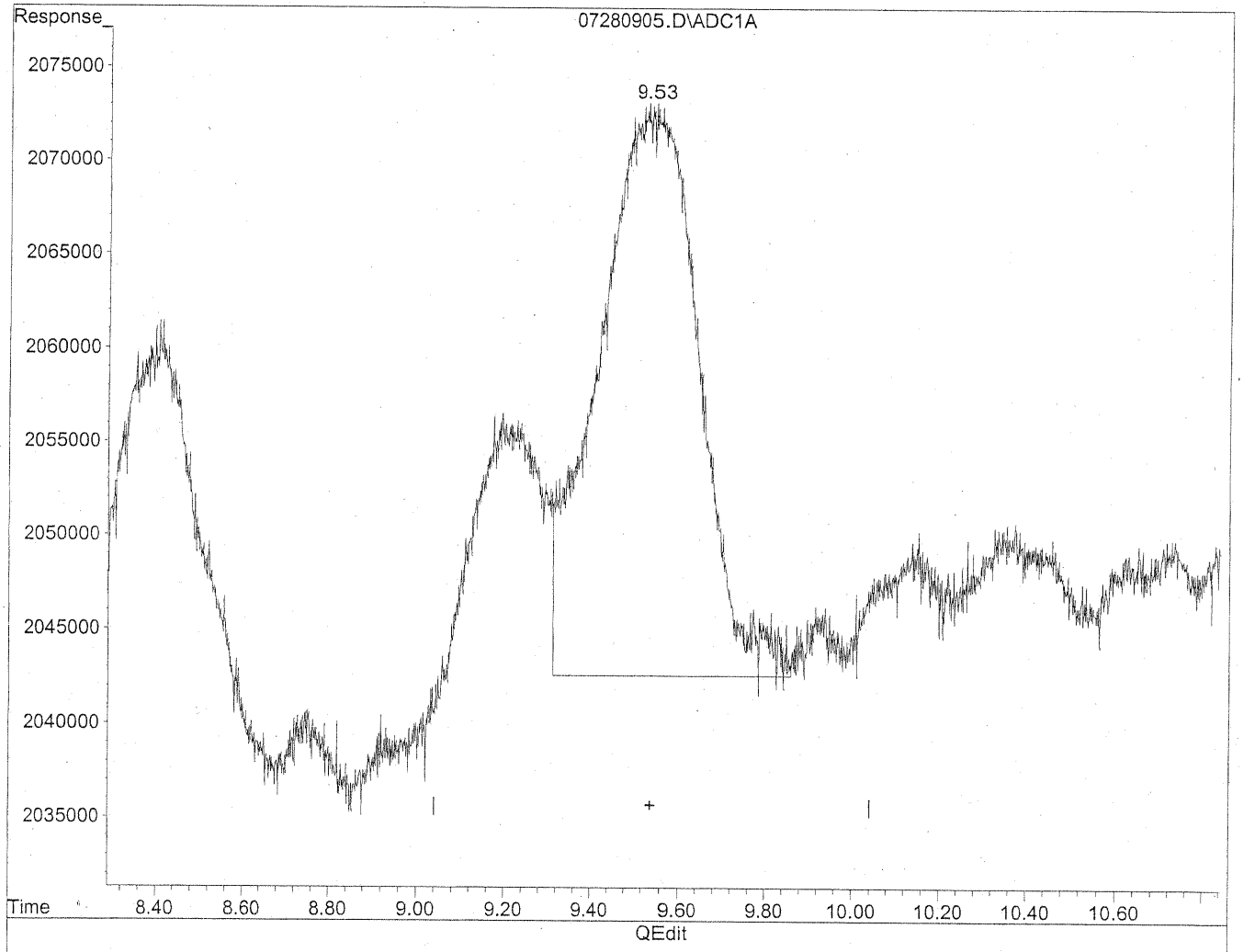
409

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280905.D Vial: 5
Acq On : 28 Jul 2009 9:39 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:25 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

9.53min 89.131ng/ml m

response 4801019

HC 7/29/09 BC

7/29/09

410

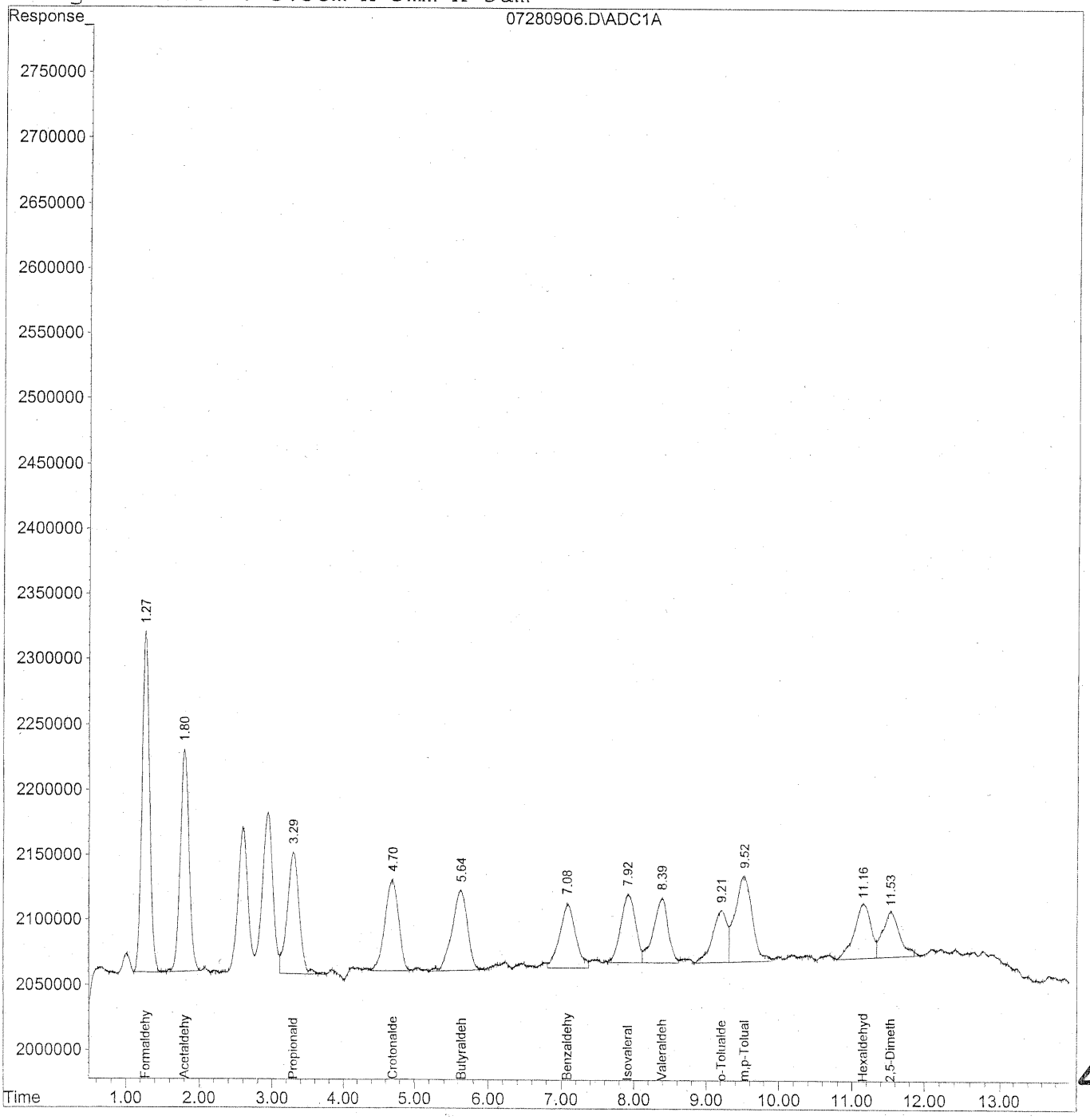
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280906.D Vial: 6
Acq On : 28 Jul 2009 9:54 am Operator: HC
Sample : 100ng/ml TO11A Std S21-07270905 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



411

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280906.D Vial: 6
Acq On : 28 Jul 2009 9:54 am Operator: HC
Sample : 100ng/ml TO11A Std S21-07270905 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

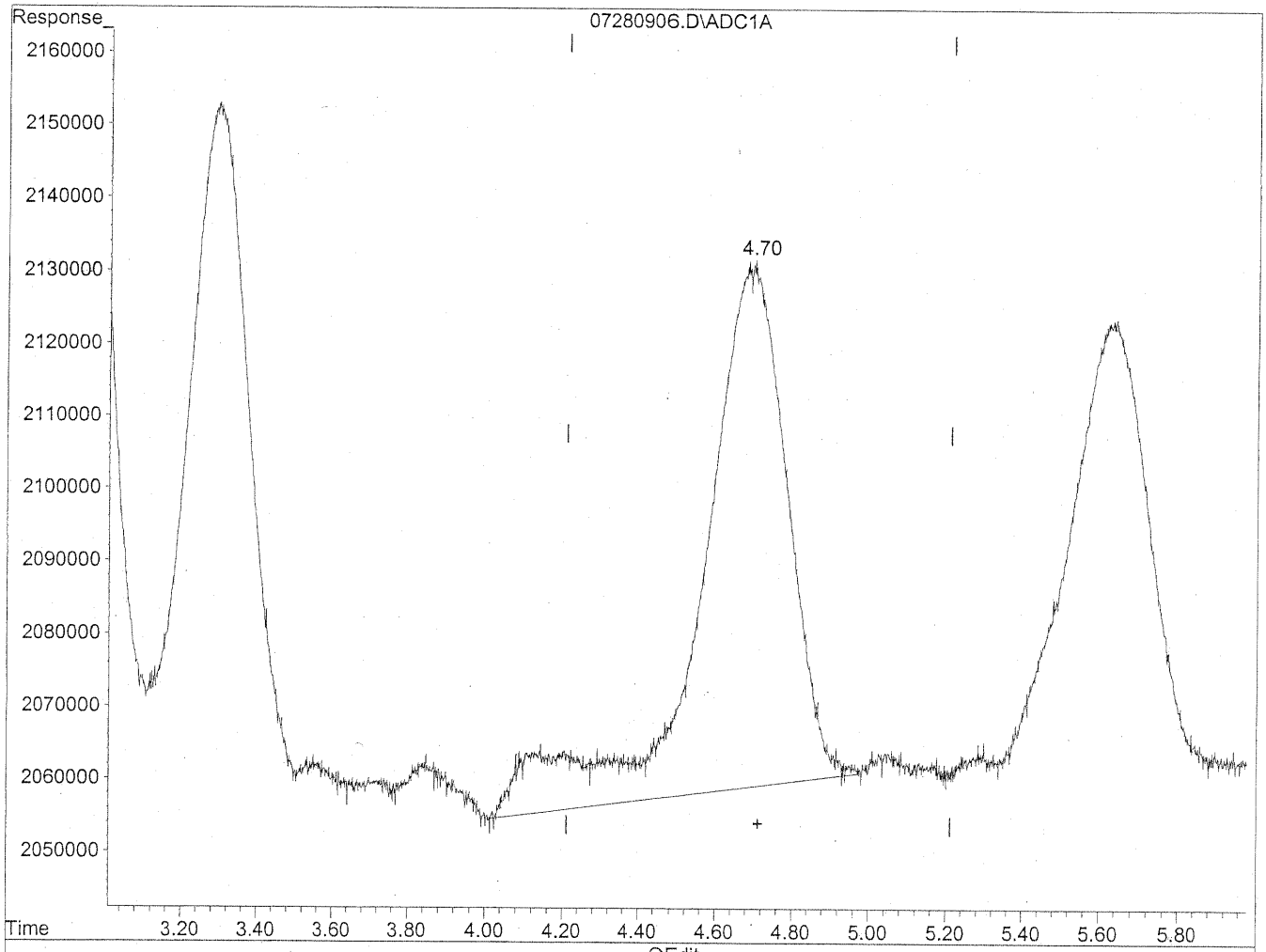
	Compound	R.T.	Response	Conc Units
	Target Compounds			
1)	Formaldehyde	1.27	18283557	104.125 ng/ml
2)	Acetaldehyde	1.80	13784712	102.185 ng/ml
3)	Propionaldehyde	3.29	10870707	105.751 ng/ml
4)	Crotonaldehyde	4.70	9346475	84.541 ng/mlm
5)	Butyraldehyde	5.63	8839595	109.842 ng/ml
6)	Benzaldehyde	7.08	7282249	115.457 ng/mlm
7)	Isovaleraldehyde	7.92	7487274	84.457 ng/ml
8)	Valeraldehyde	8.39	7060988	84.975 ng/ml
9)	o-Tolualdehyde	9.21	5548699	103.001 ng/ml
10)	m,p-Tolualdehyde	9.52	10979457	203.834 ng/ml
11)	Hexaldehyde	11.16	6702769	99.835 ng/ml
12)	2,5-Dimethylbenzaldehyde	11.53	5798505	111.652 ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280906.D Vial: 6
Acq On : 28 Jul 2009 9:54 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:29 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.69min 102.369ng/ml

response 11317409

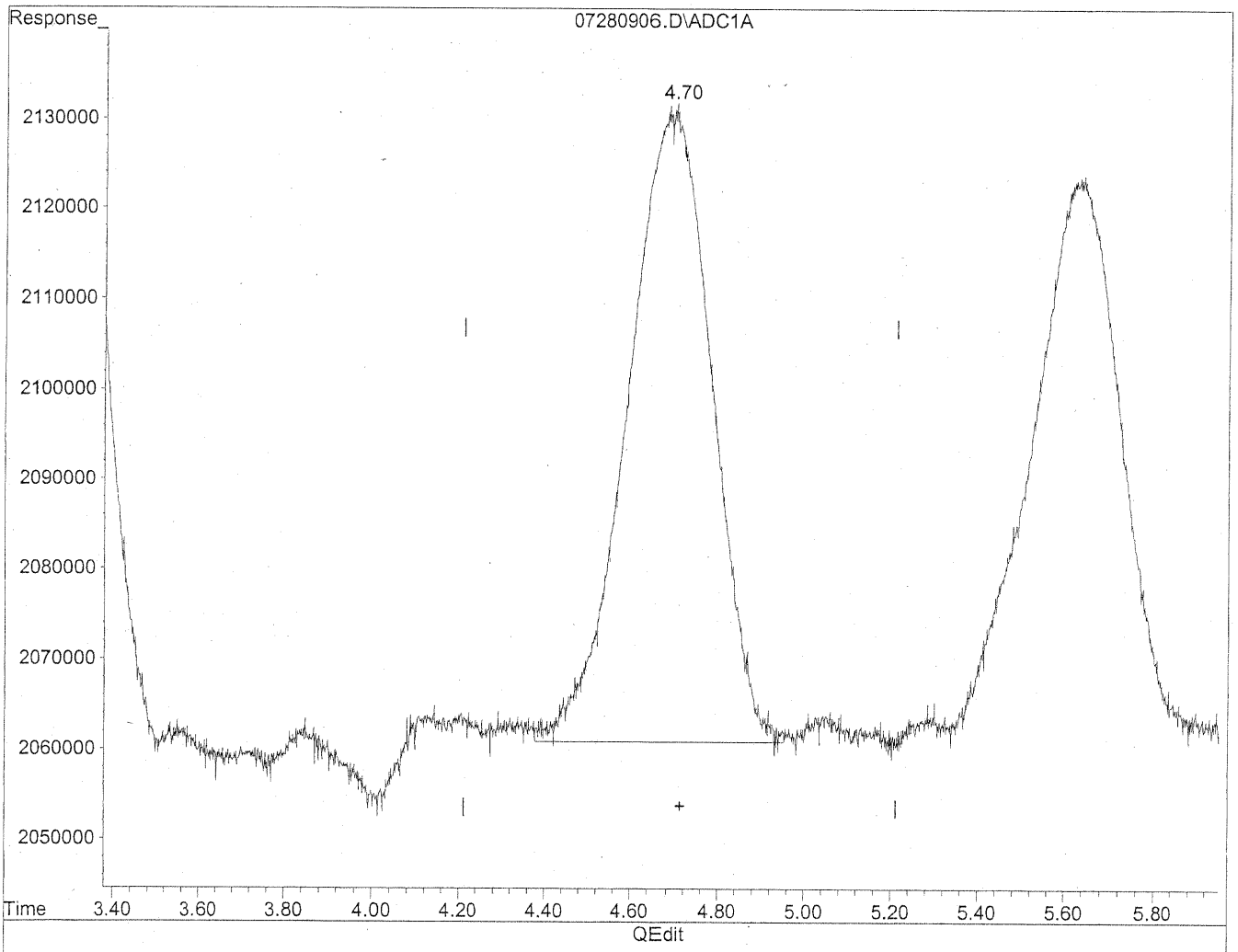
413

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280906.D Vial: 6
Acq On : 28 Jul 2009 9:54 am Operator: HC
Sample : 100ng/ml TO11A Std S21-07270905 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 15:29:52 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.70min 84.541ng/ml m

response 9346475

HC 7/28/09

KK 7/29/09

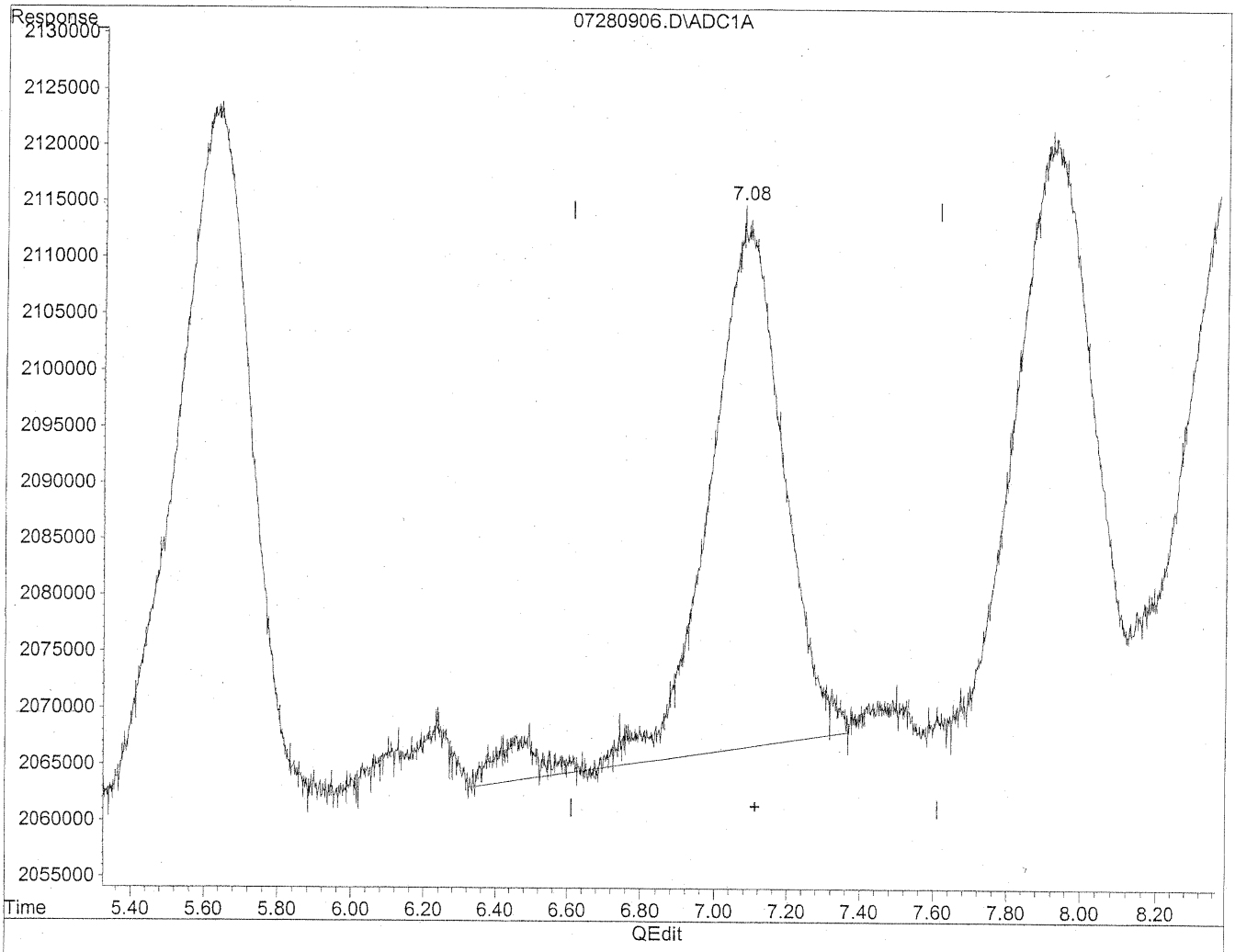
414

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280906.D Vial: 6
Acq On : 28 Jul 2009 9:54 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:29 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

7.09min 108.123ng/ml

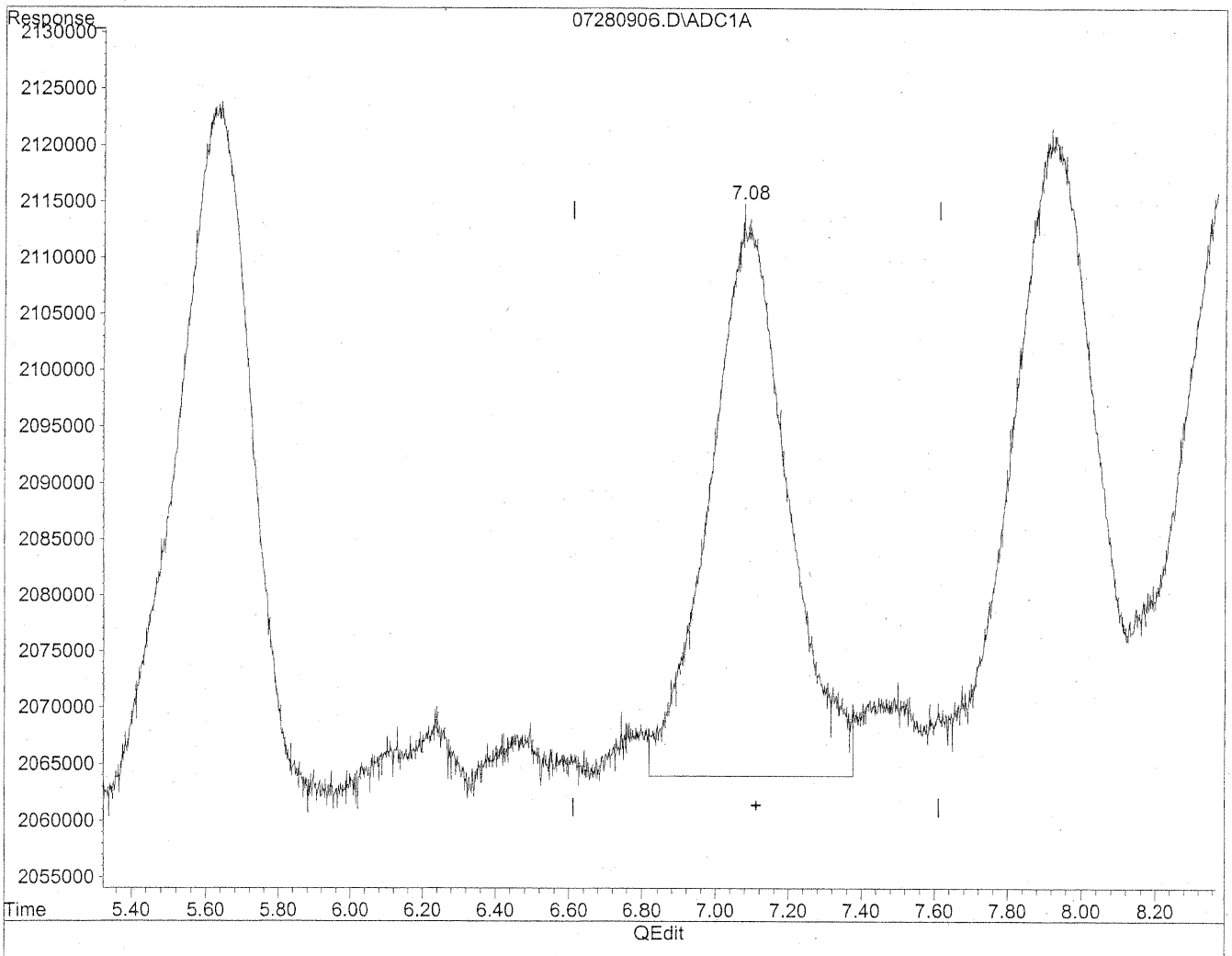
response 6819663

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280906.D Vial: 6
Acq On : 28 Jul 2009 9:54 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:29 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

7.08min 115.457ng/ml m

response 7282249

tc 7/28/09 JC

KR 7/29/09

(+) = Expected Retention Time

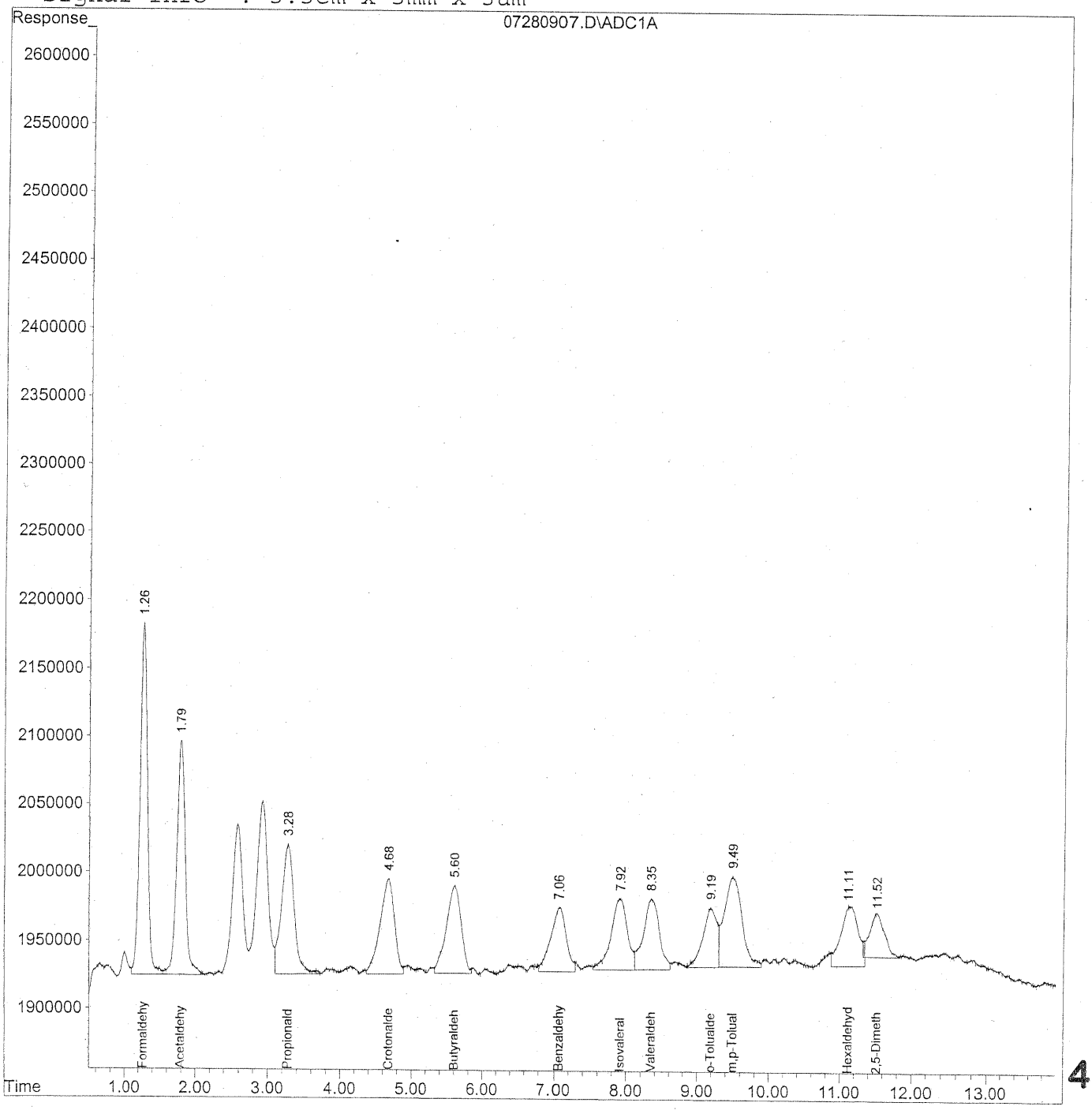
07280906.D TO110709.M Tue Jul 28 14:30:30 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std S21-07270905 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:33 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



417

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std S21-07270905 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:33 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

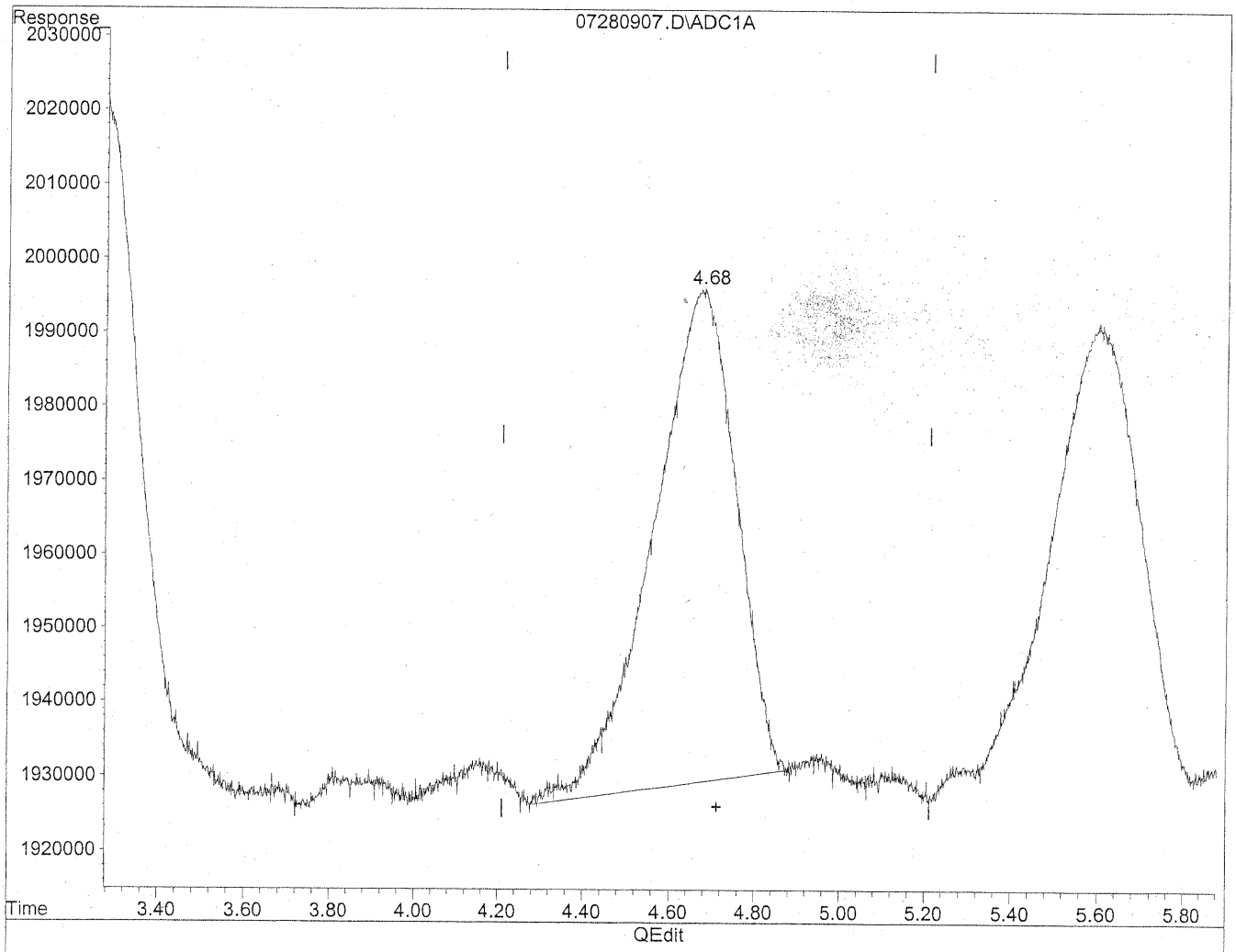
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.26	18449443	105.069	ng/ml
2)	Acetaldehyde	1.79	14434553	107.002	ng/ml
3)	Propionaldehyde	3.28	11389784	110.800	ng/ml
4)	Crotonaldehyde	4.68	9814490	88.774	ng/mlm
5)	Butyraldehyde	5.60	9432197	117.206	ng/mlm
6)	Benzaldehyde	7.06	6706722	106.332	ng/mlm
7)	Isovaleraldehyde	7.92	8338385	94.058	ng/mlm
8)	Valeraldehyde	8.35	8117341	97.688	ng/mlm
9)	o-Tolualdehyde	9.19	5921917	109.929	ng/mlm
10)	m,p-Tolualdehyde	9.49	11235135	208.581	ng/mlm
11)	Hexaldehyde	11.11	7714022	114.897	ng/mlm
12)	2,5-Dimethylbenzaldehyde	11.51	4735227	91.178	ng/mlm

(f)=RT Delta > 1/2 Window (m)=manual int.
07280907.D TO110709.M Tue Jul 28 17:19:21 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.67min 80.883ng/ml

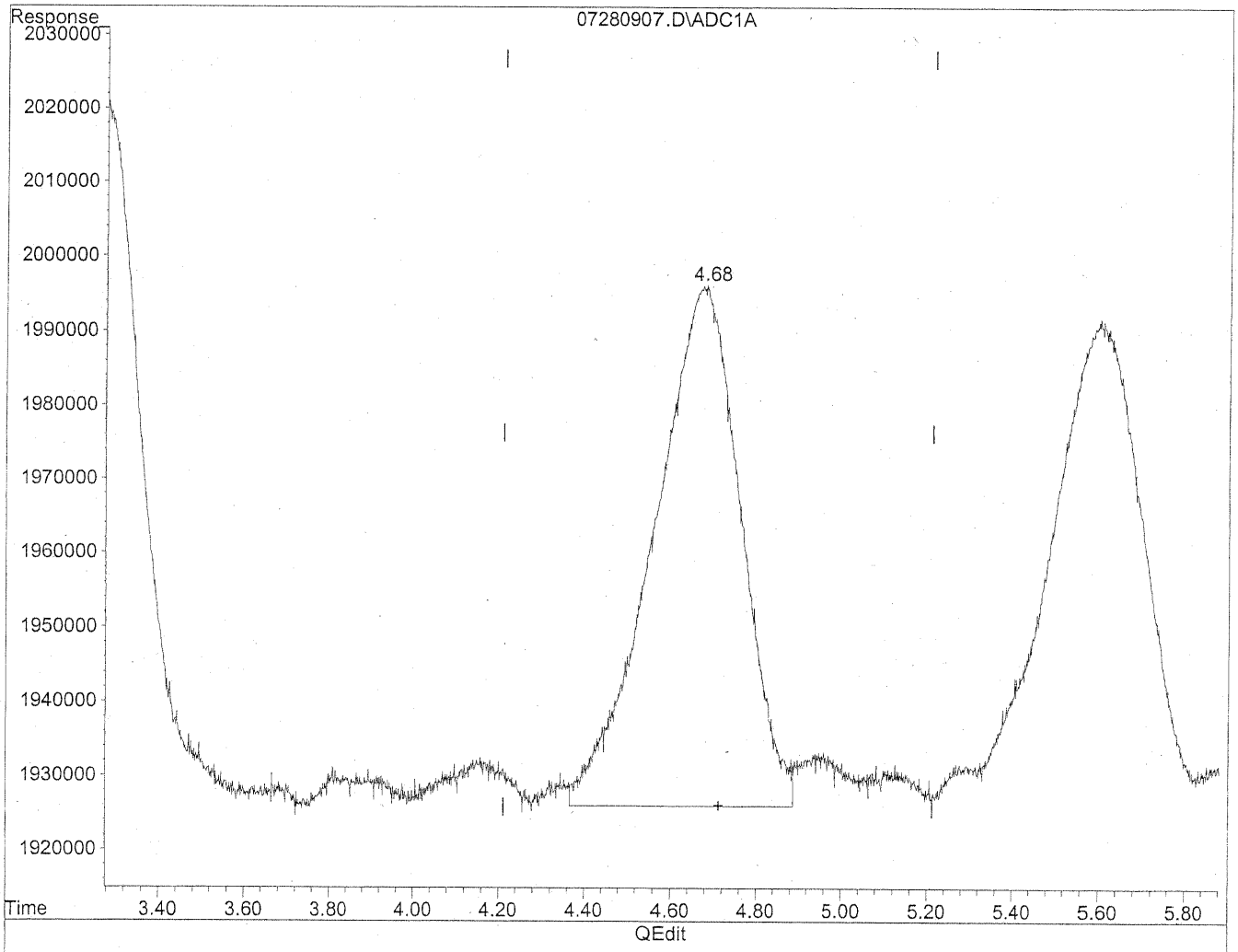
response 8942013

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.68min 88.774ng/ml m

response 9814490

AC 7/28/09 1C

KR 7/29/09

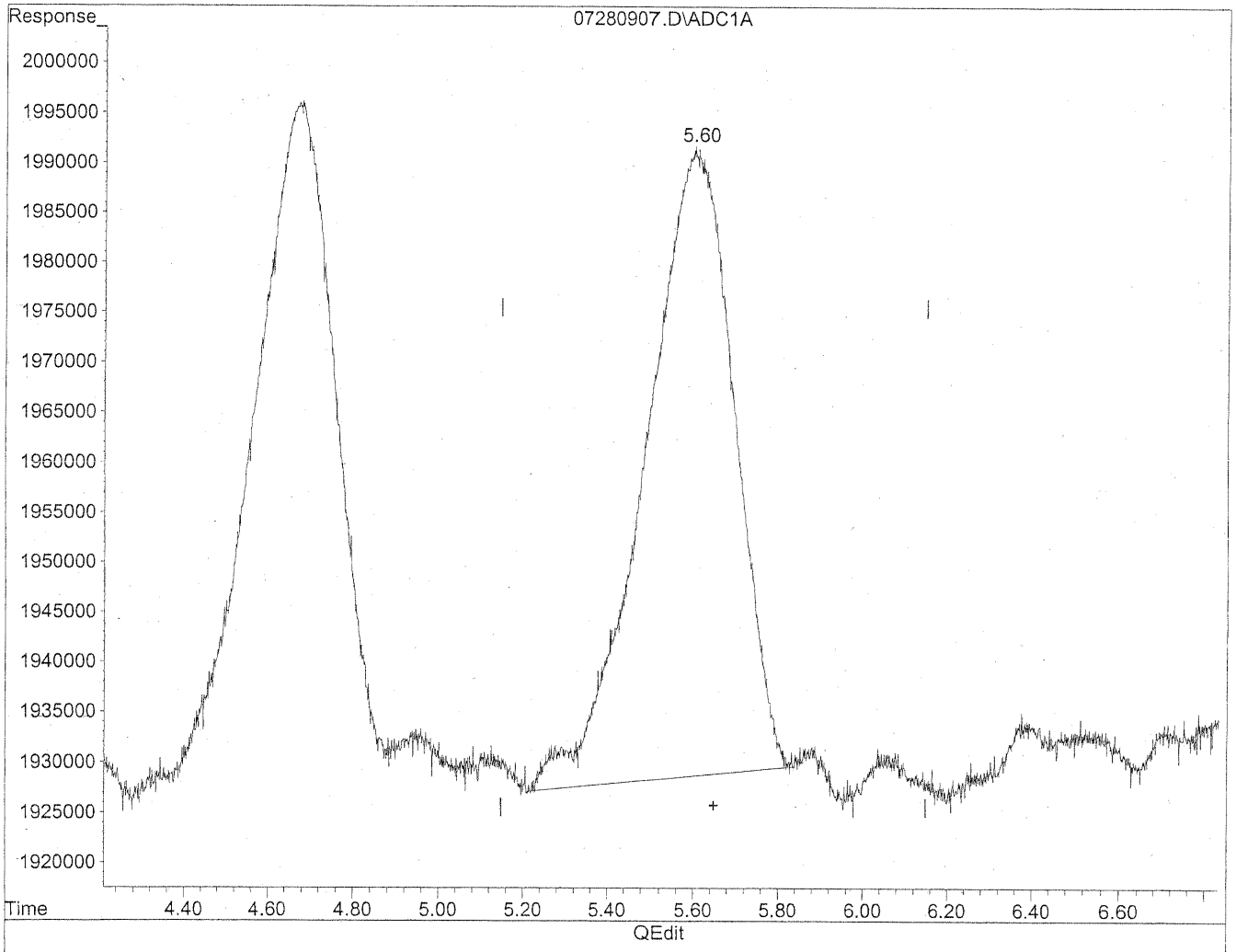
(+) = Expected Retention Time

07280907.D TO110709.M Tue Jul 28 14:31:09 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.60min 112.634ng/ml

response 9064274

421

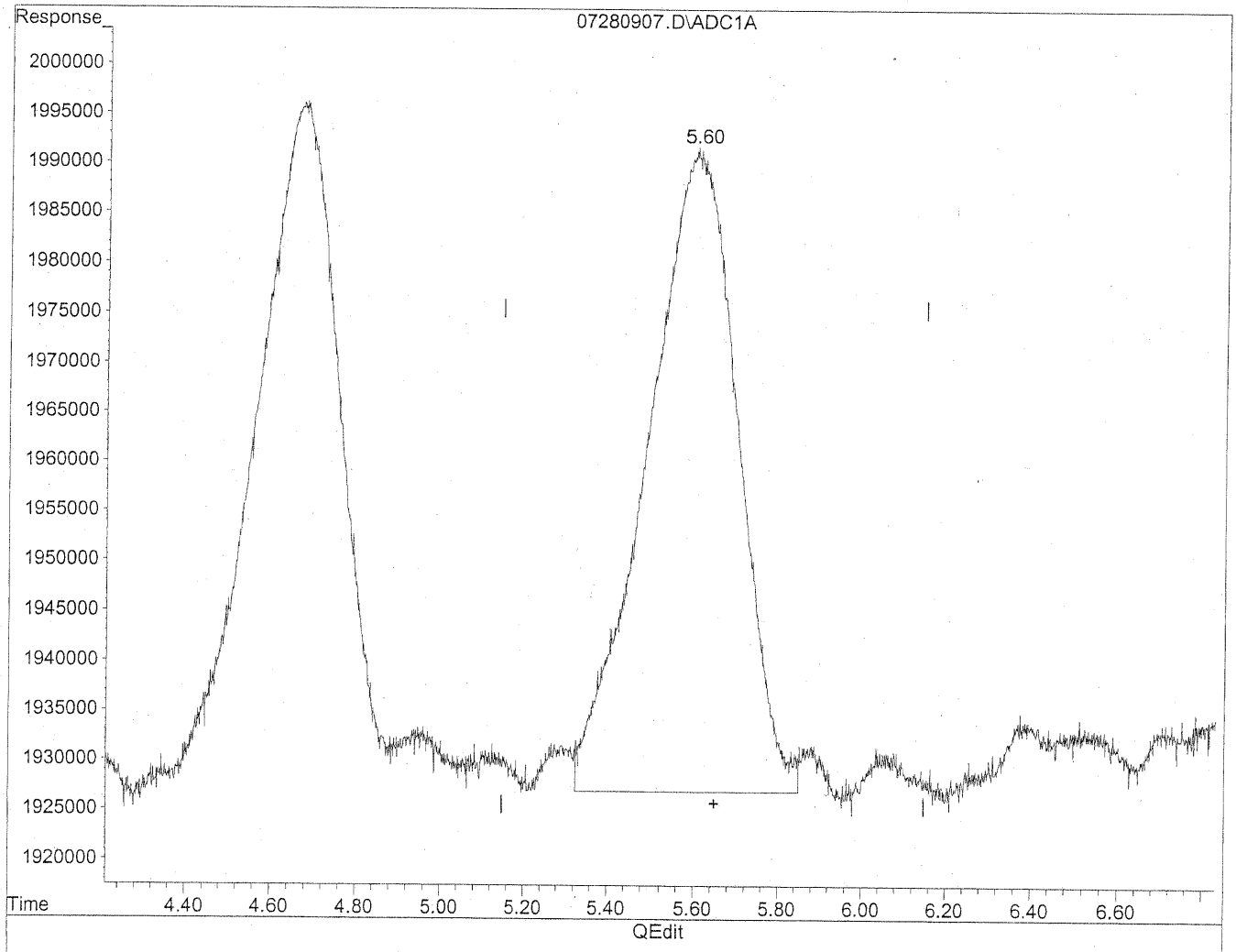
(+) = Expected Retention Time

07280907.D TO110709.M Tue Jul 28 14:31:17 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.60min 117.206ng/ml m

response 9432197

HC station 1C

7/29/09

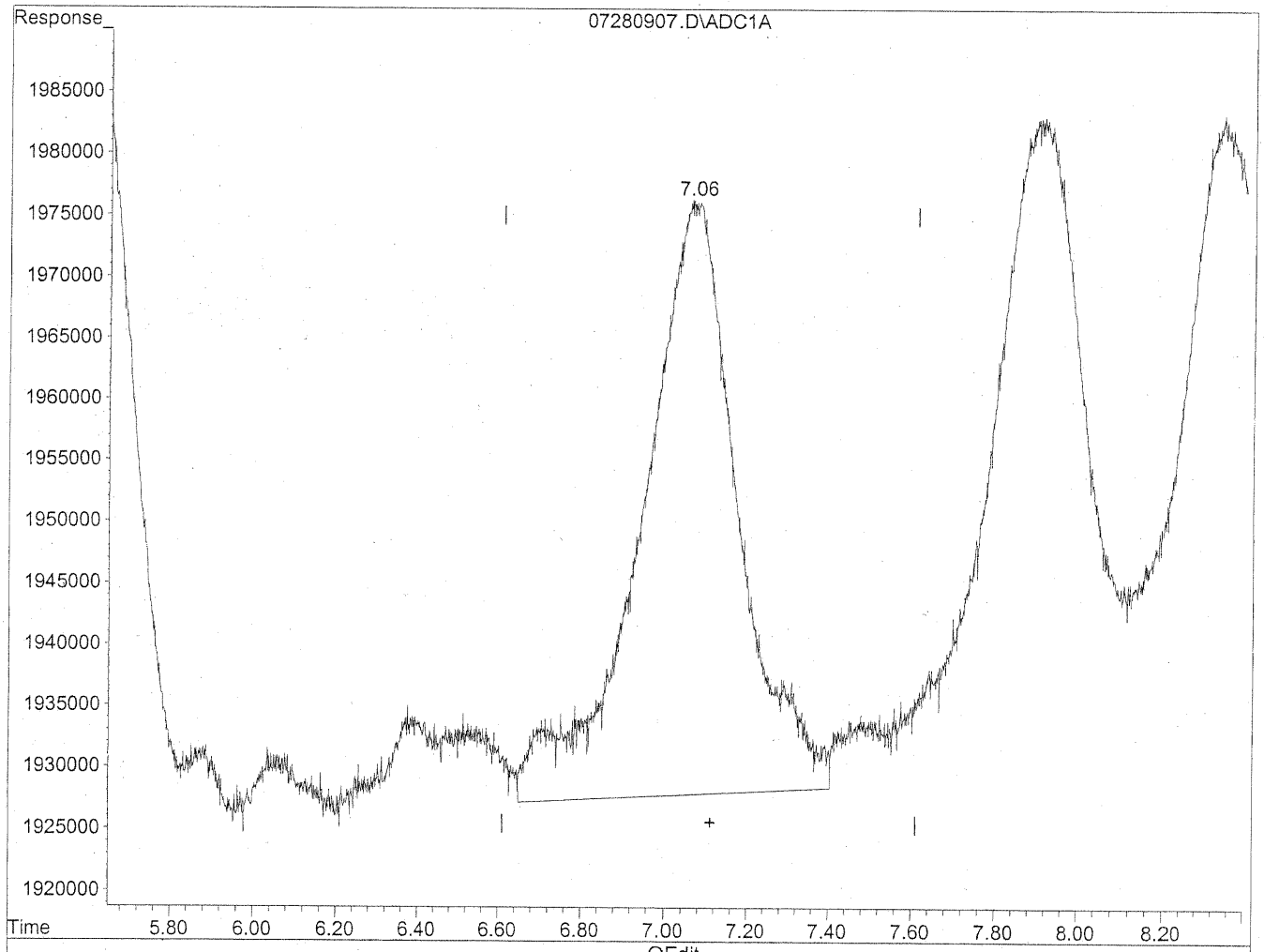
422

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

7.07min 123.223ng/ml

response 7772036

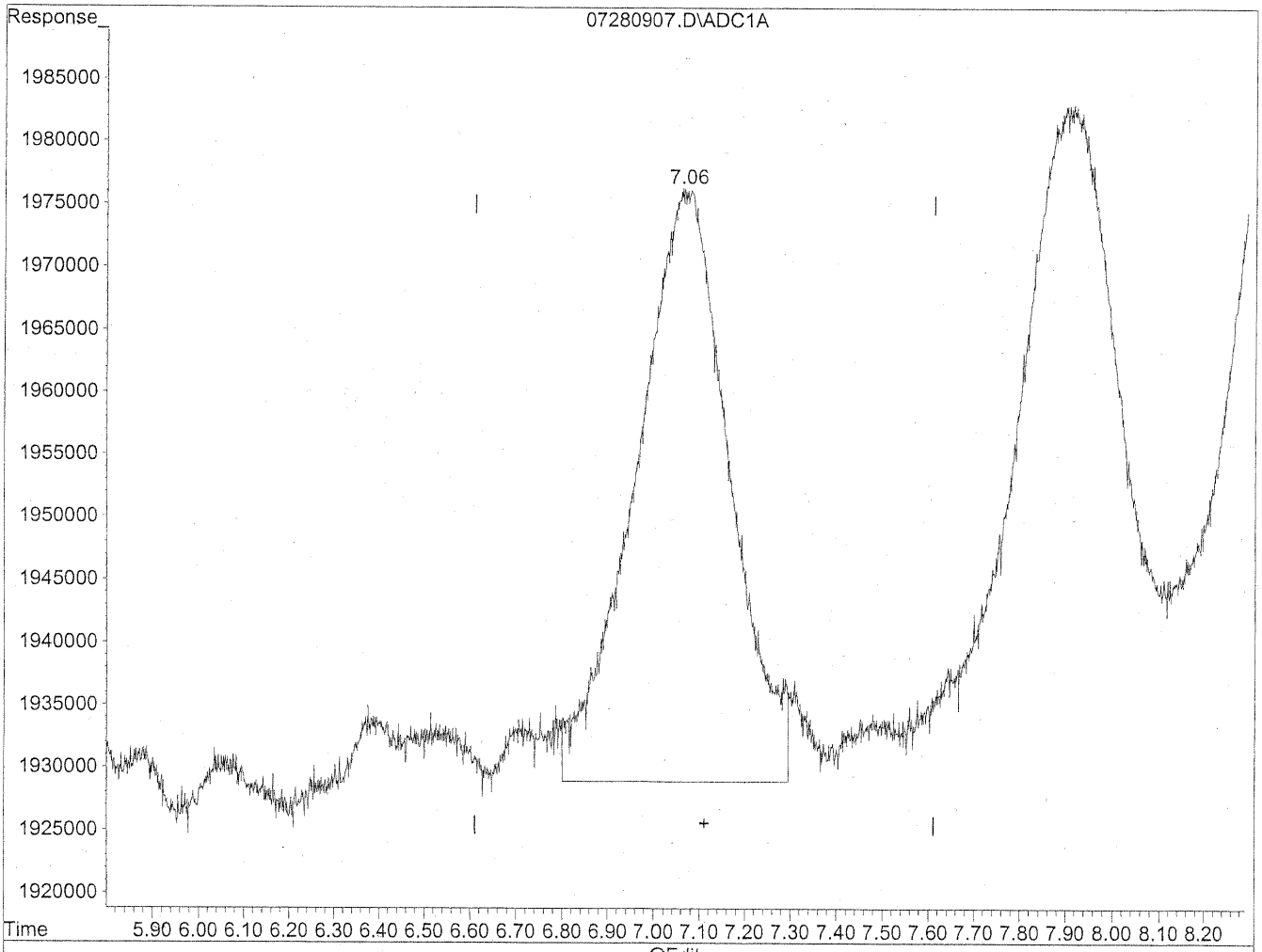
(+) = Expected Retention Time

07280907.D TO110709.M Tue Jul 28 14:31:29 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

7.06min 106.332ng/ml m

response 6706722

HC 7/28/09 LC

7/24/09

424

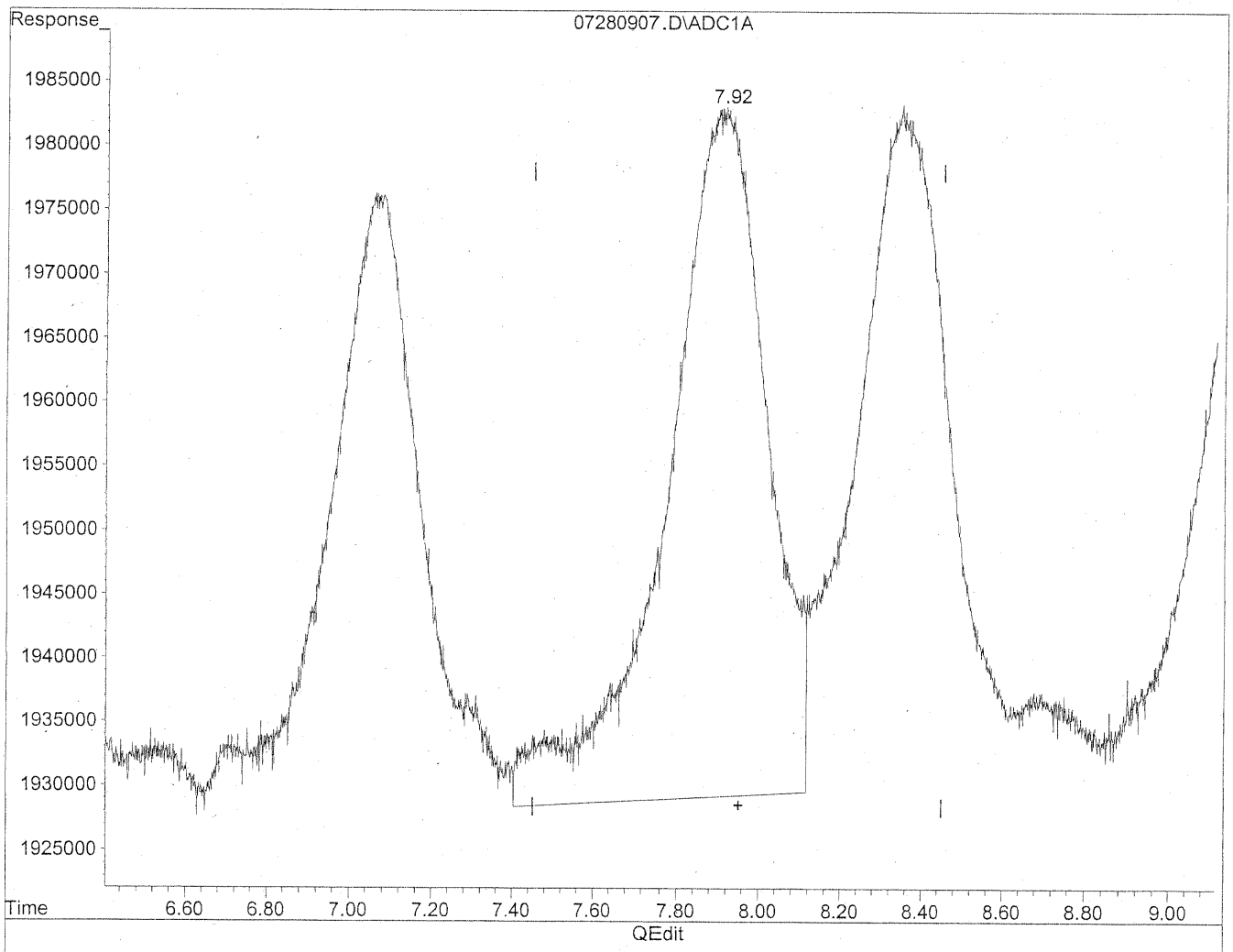
(+) = Expected Retention Time

07280907.D TO110709.M Tue Jul 28 14:31:42 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.91min 103.108ng/ml

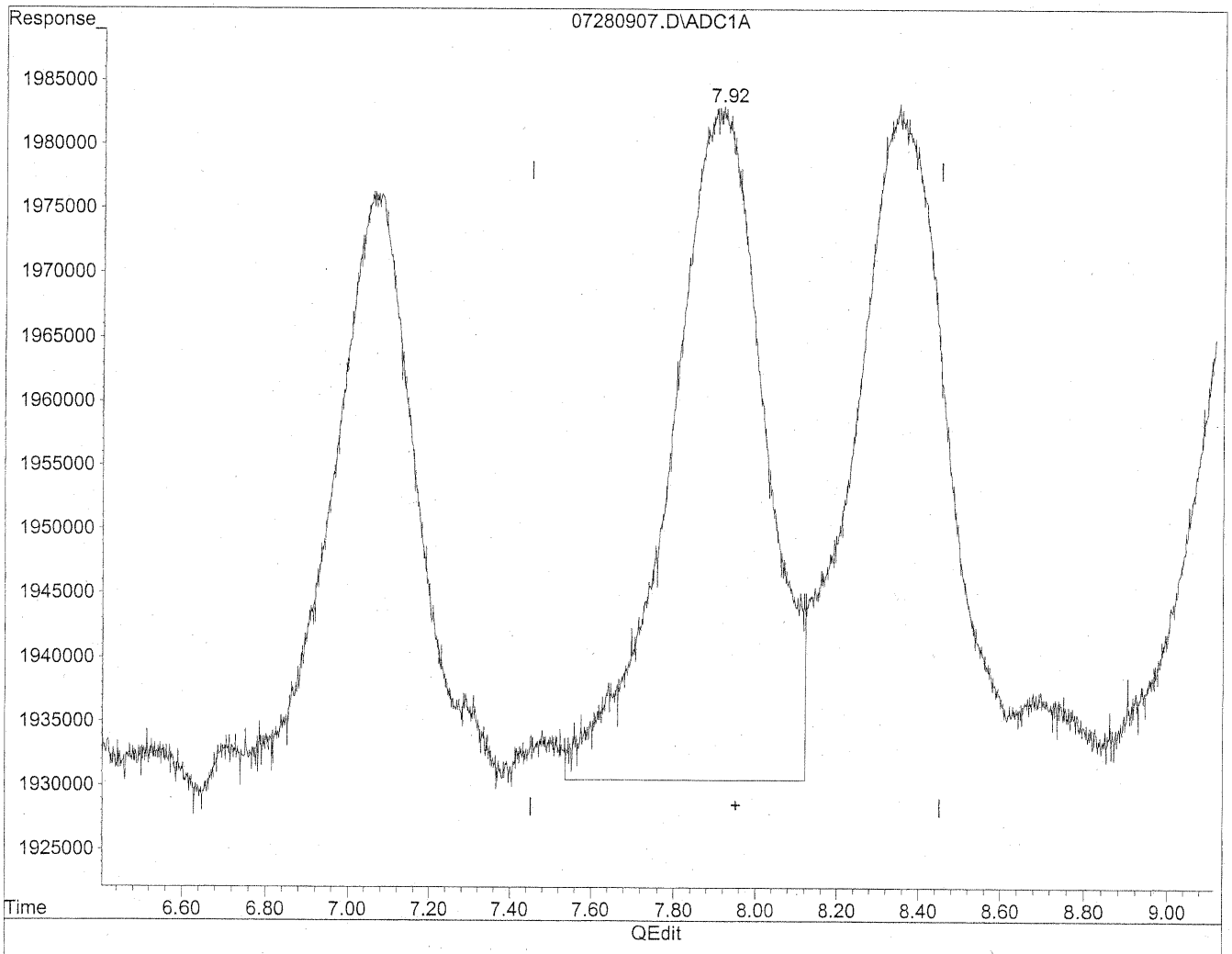
response 9140643

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.92min 94.058ng/ml m

response 8338385

7/28/09 IC

7/29/09

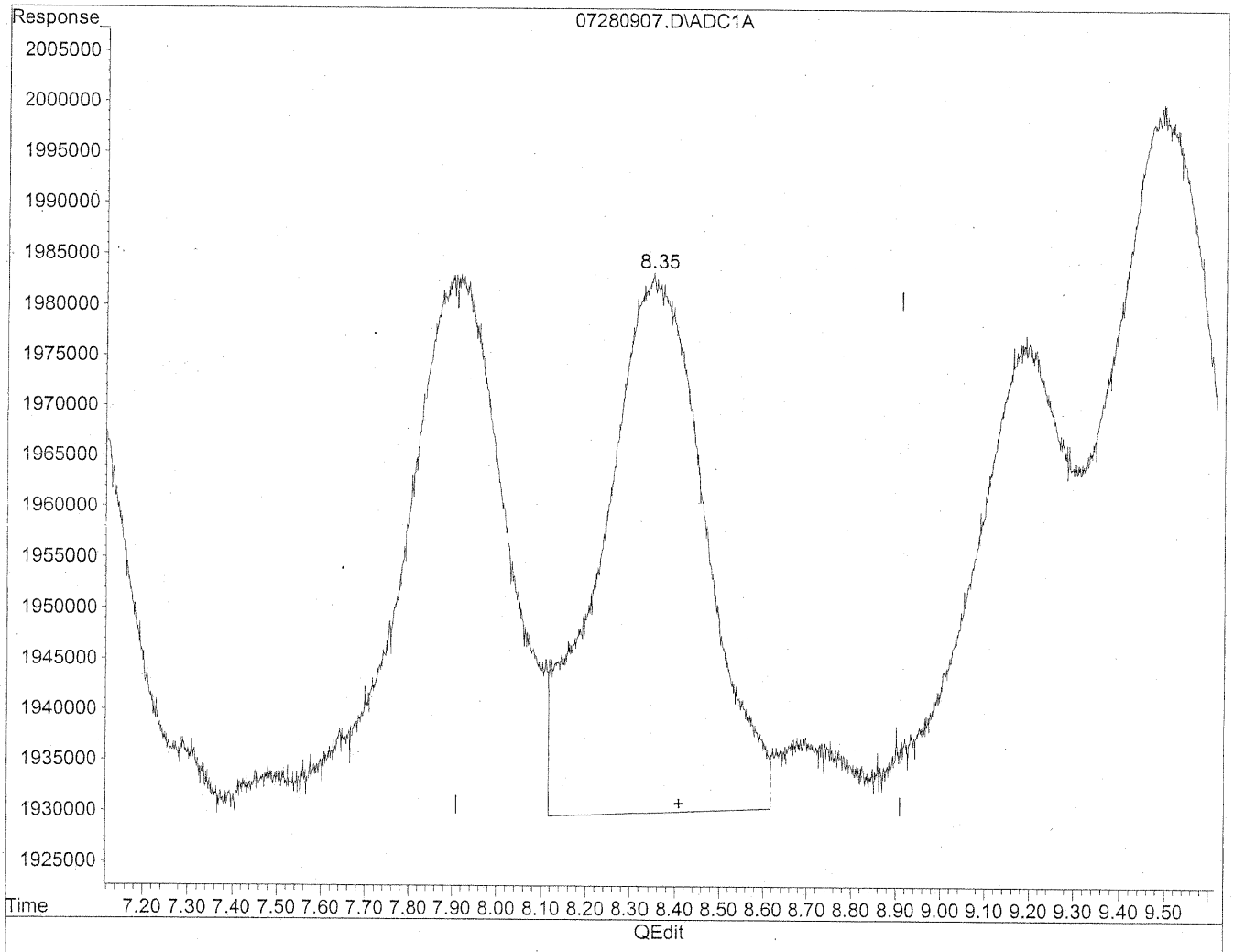
426

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

8.35min 101.373ng/ml

response 8423554

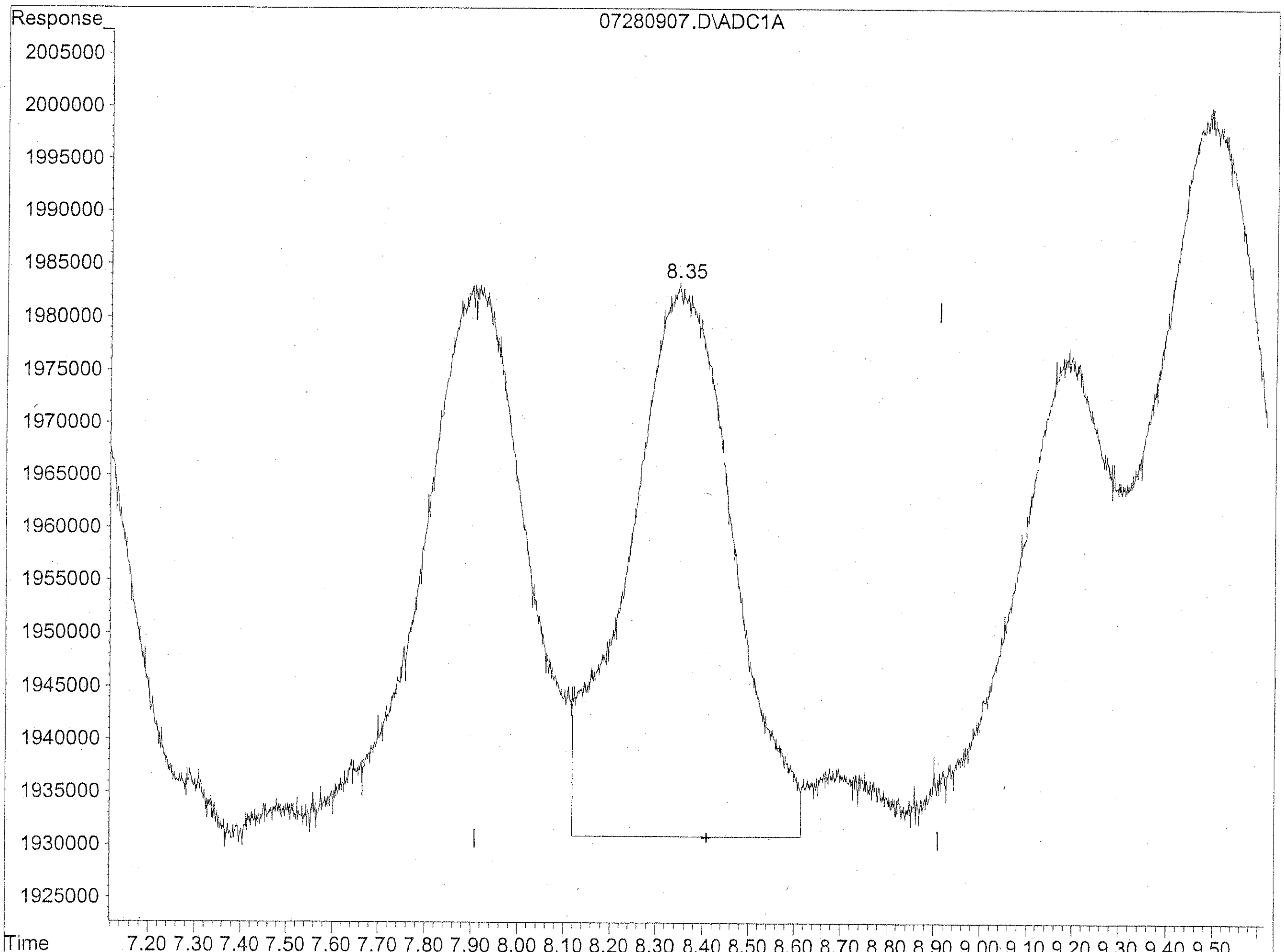
427

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D   Vial: 7
Acq On : 28 Jul 2009 10:09 am   Operator: HC
Sample : 100ng/ml TO11A Std   Inst : LC 01
Misc :   Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109   Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



QEdit

(8) Valeraldehyde

8.35min 97.688ng/ml m

response 8117341

HC 7/28/09 BC

KR 7/29/09

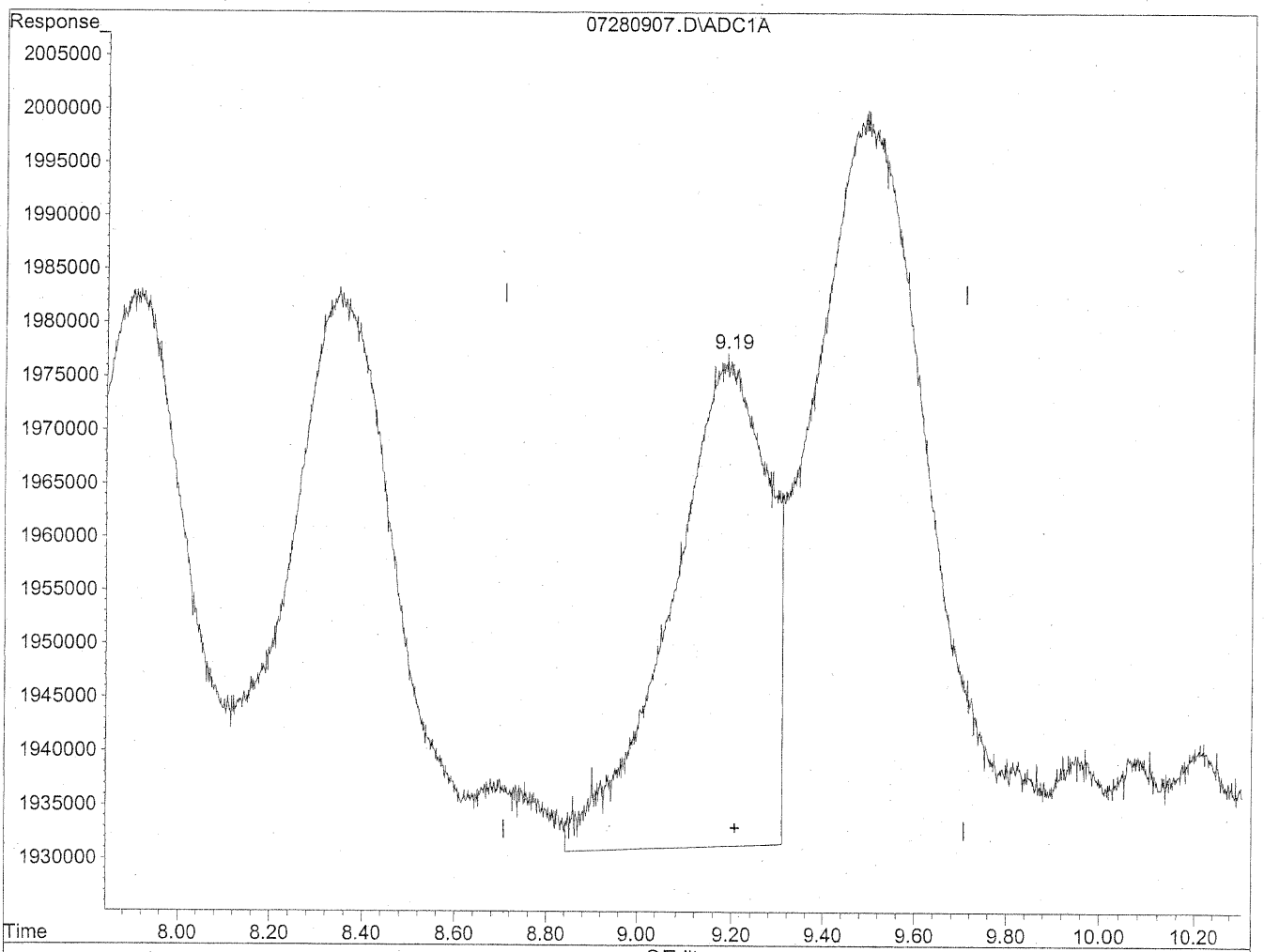
(+) = Expected Retention Time

07280907.D TO110709.M Tue Jul 28 14:32:16 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.19min 121.312ng/ml

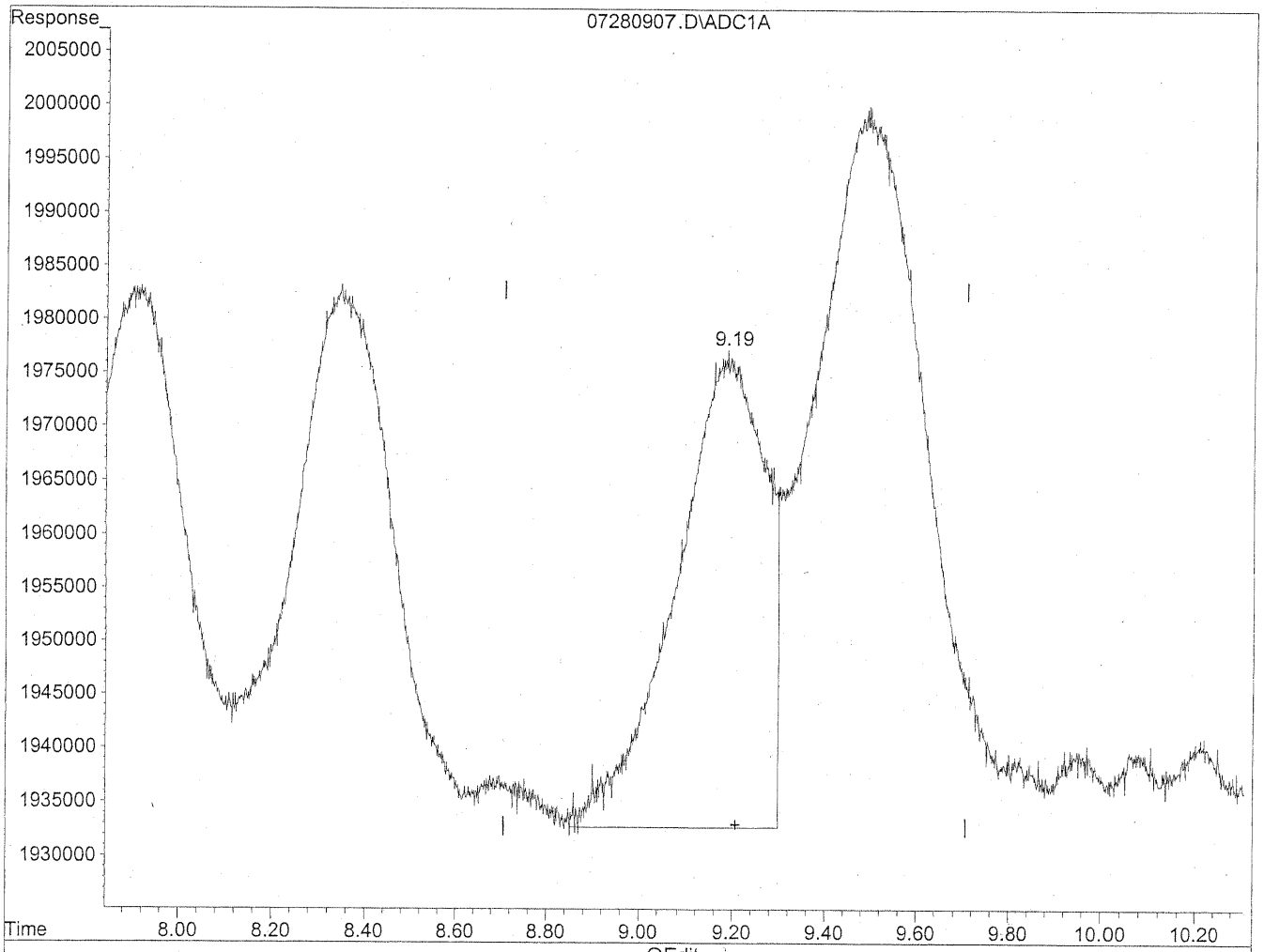
response 6535124

(+) = Expected Retention Time
07280907.D TO110709.M Tue Jul 28 14:32:23 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.19min 109.929ng/ml m

response 5921917

HK 7/28/09 BC

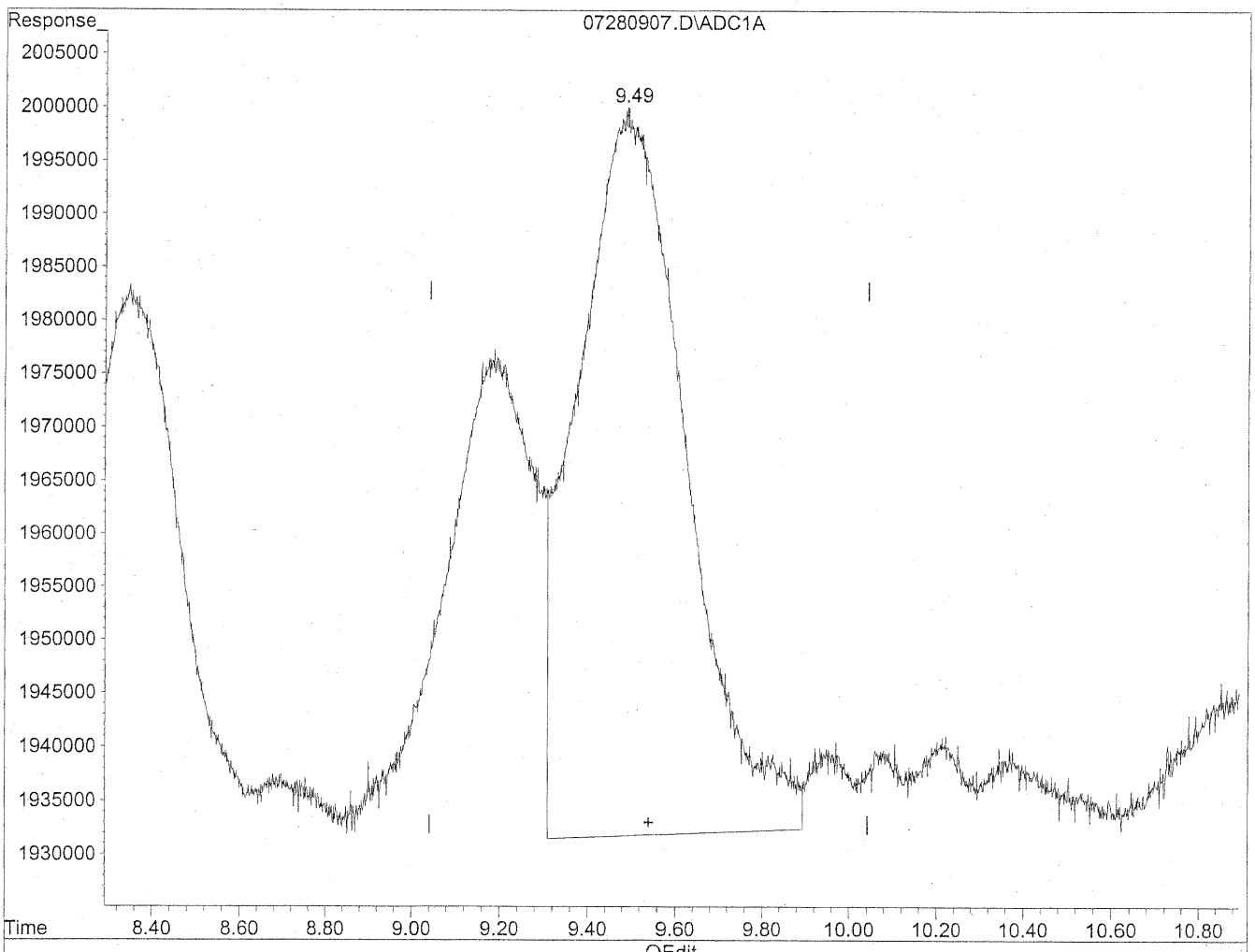
KR 7/29/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

9.49min 217.917ng/ml

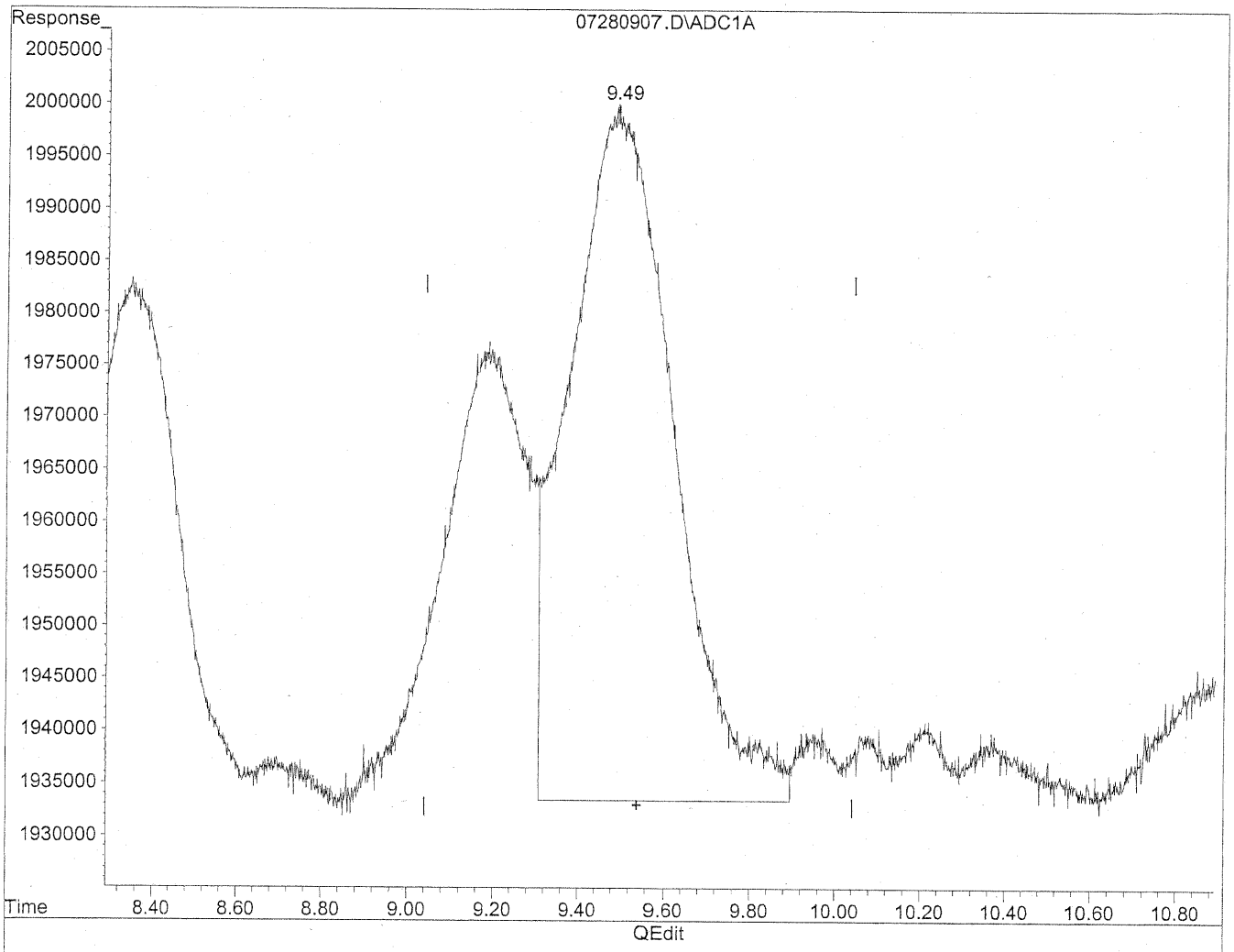
response 11738041

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

9.49min 208.581ng/ml m

response 11235135

HC 7/28/09 TBC

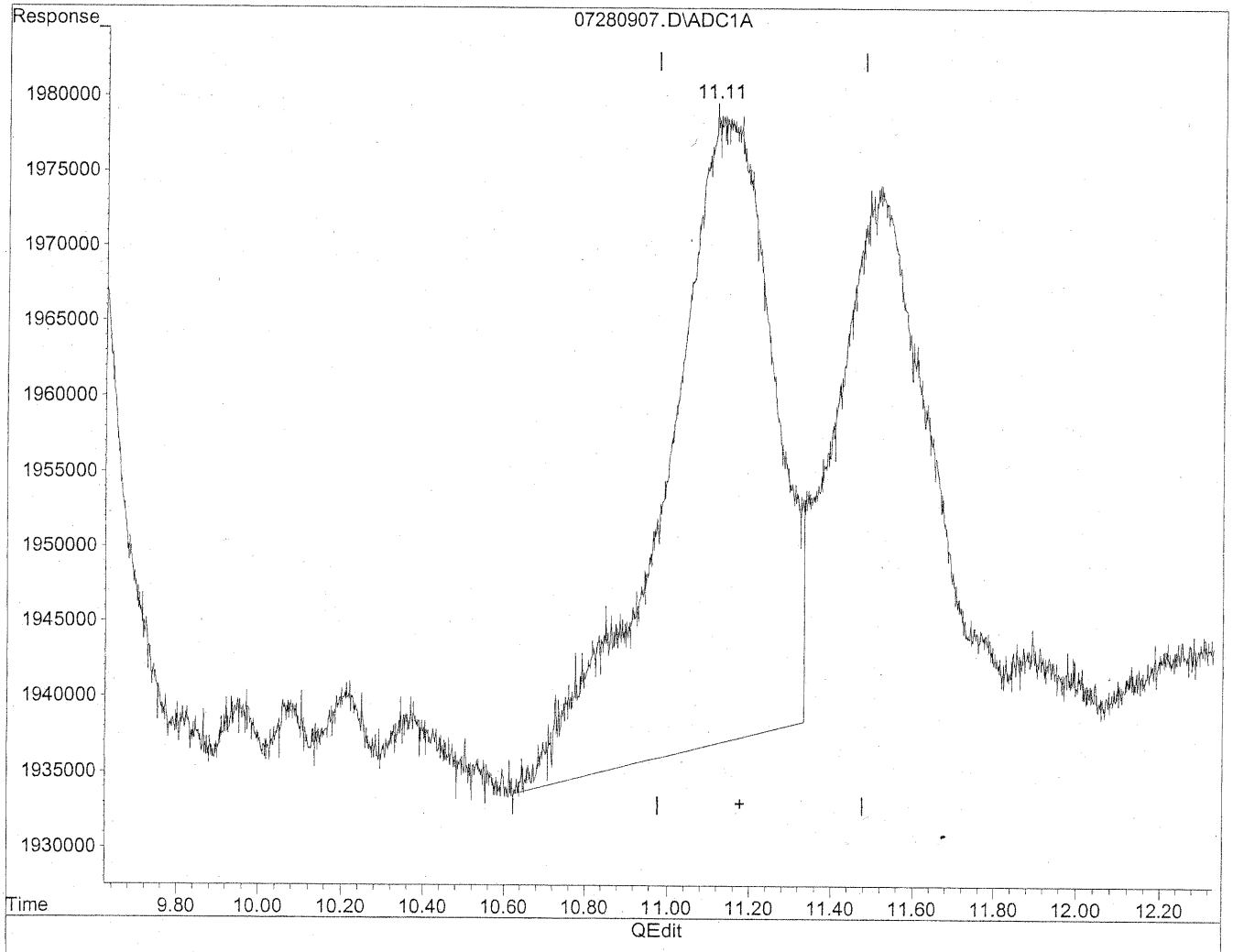
KR 7/29/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

11.14min 112.492ng/ml

response 7552544

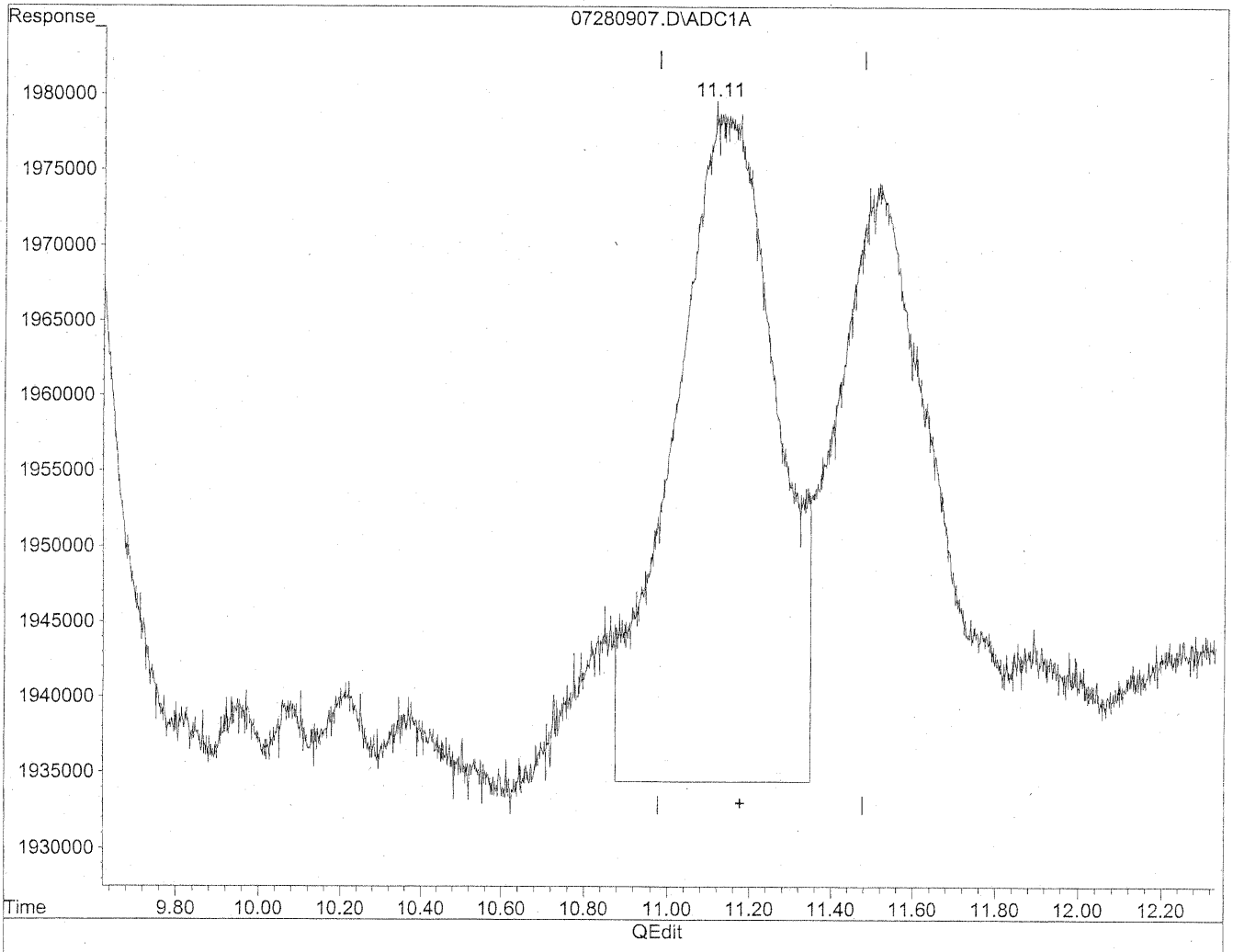
(+) = Expected Retention Time

07280907.D TO110709.M Tue Jul 28 14:32:57 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

11.11min 114.897ng/ml m

response 7714022

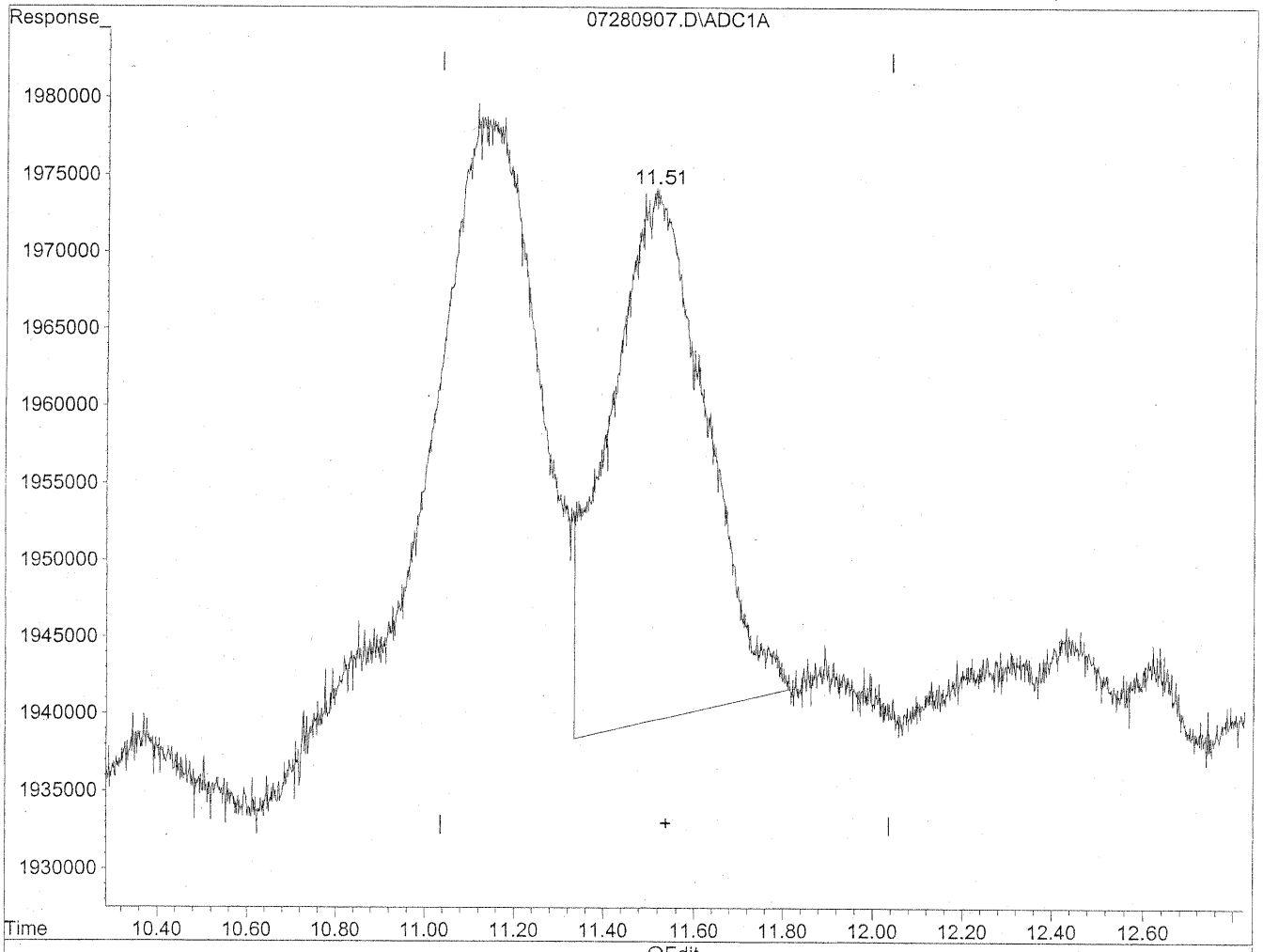
434

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

11.52min 97.911ng/ml

response 5084888

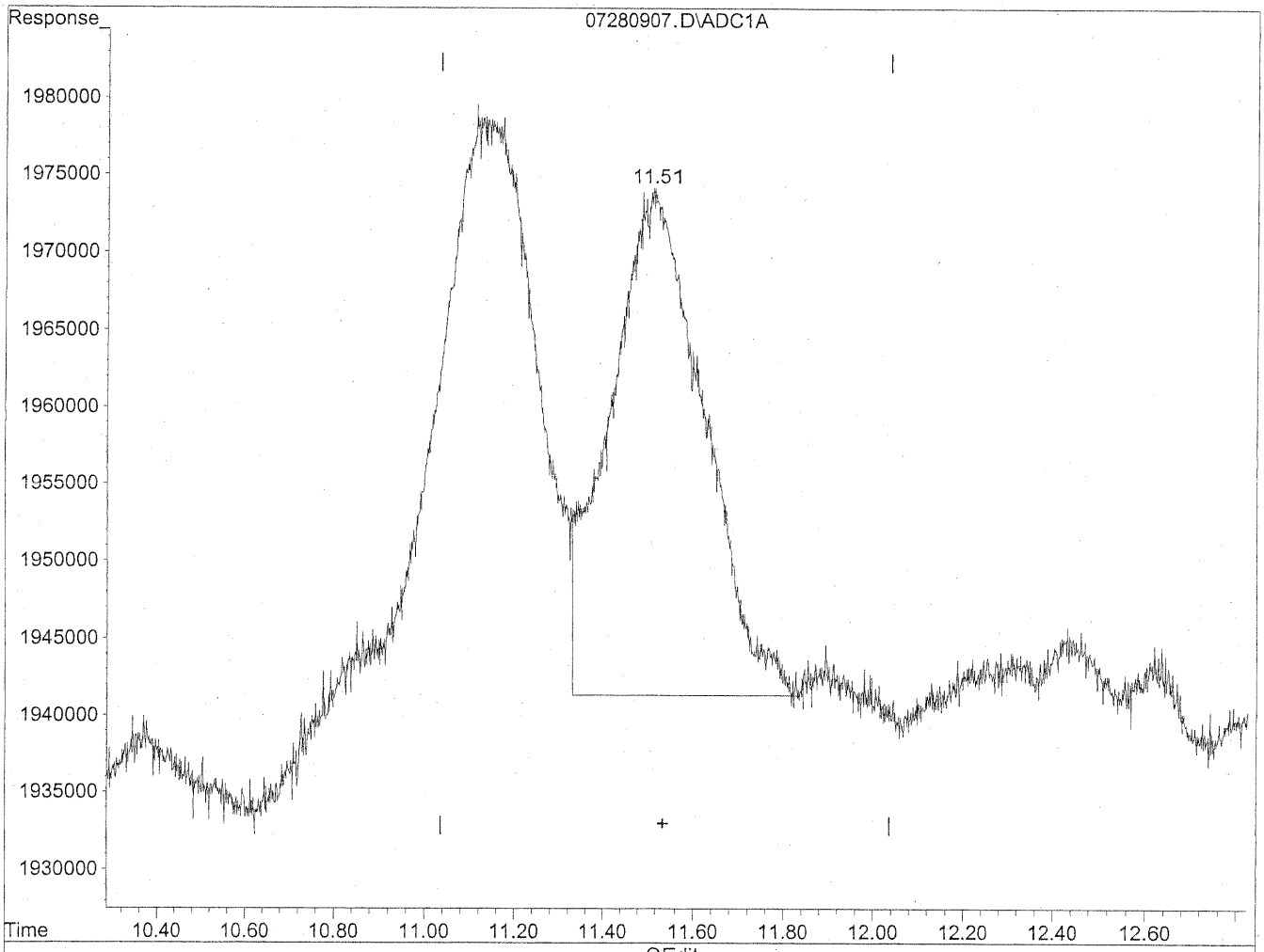
435

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



QEdit

(12) 2,5-Dimethylbenzaldehyde

11.51min 91.178ng/ml m

response 4735227

HL 7/28/09 PL

KK 7/29/09

436

(+) = Expected Retention Time

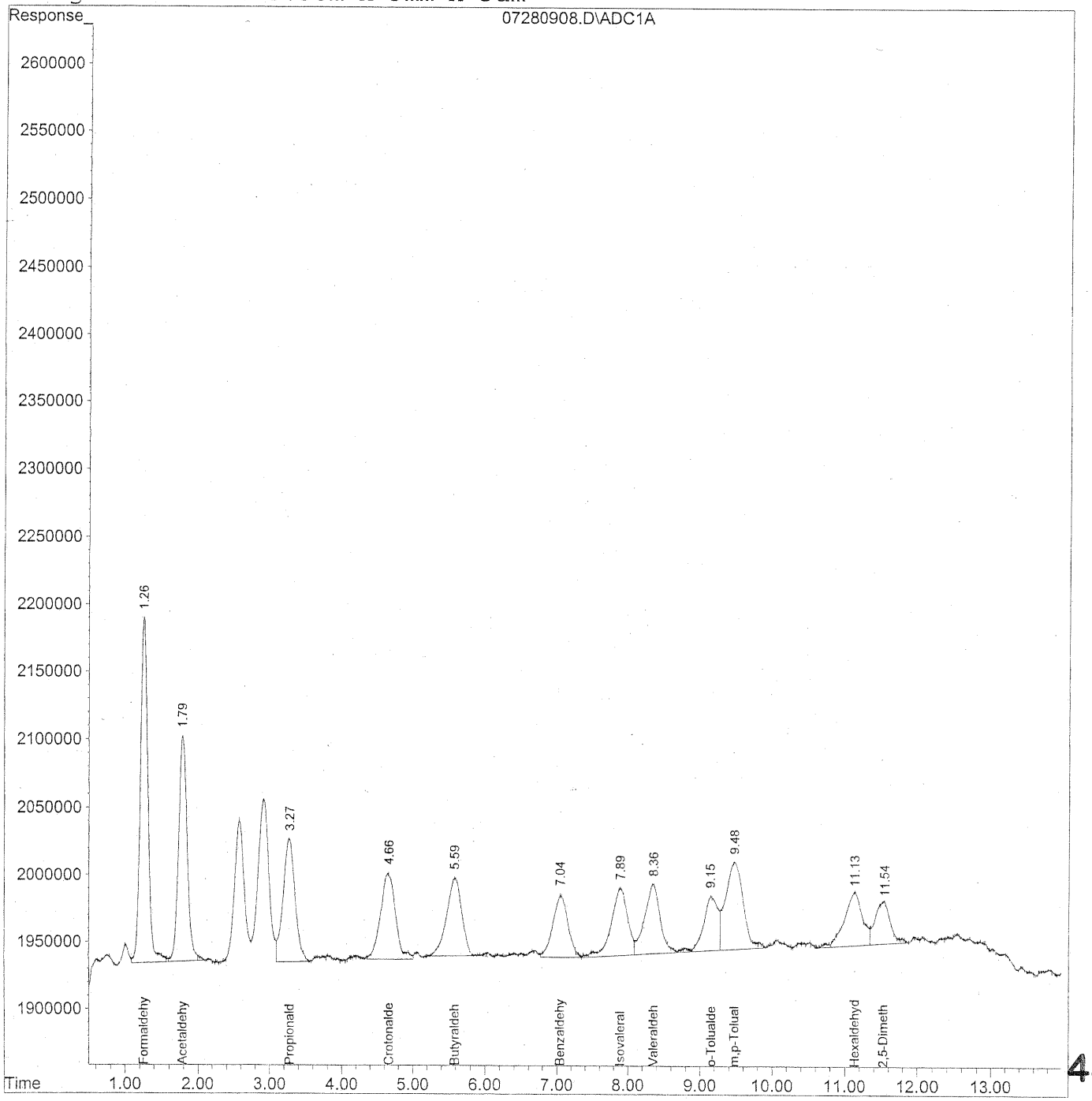
07280907.D TO110709.M Tue Jul 28 14:33:31 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280908.D Vial: 8
Acq On : 28 Jul 2009 10:24 am Operator: HC
Sample : 100ng/ml TO11A Std S21-07270905 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:34 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



437

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280908.D Vial: 8
Acq On : 28 Jul 2009 10:24 am Operator: HC
Sample : 100ng/ml TO11A Std S21-07270905 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:34 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

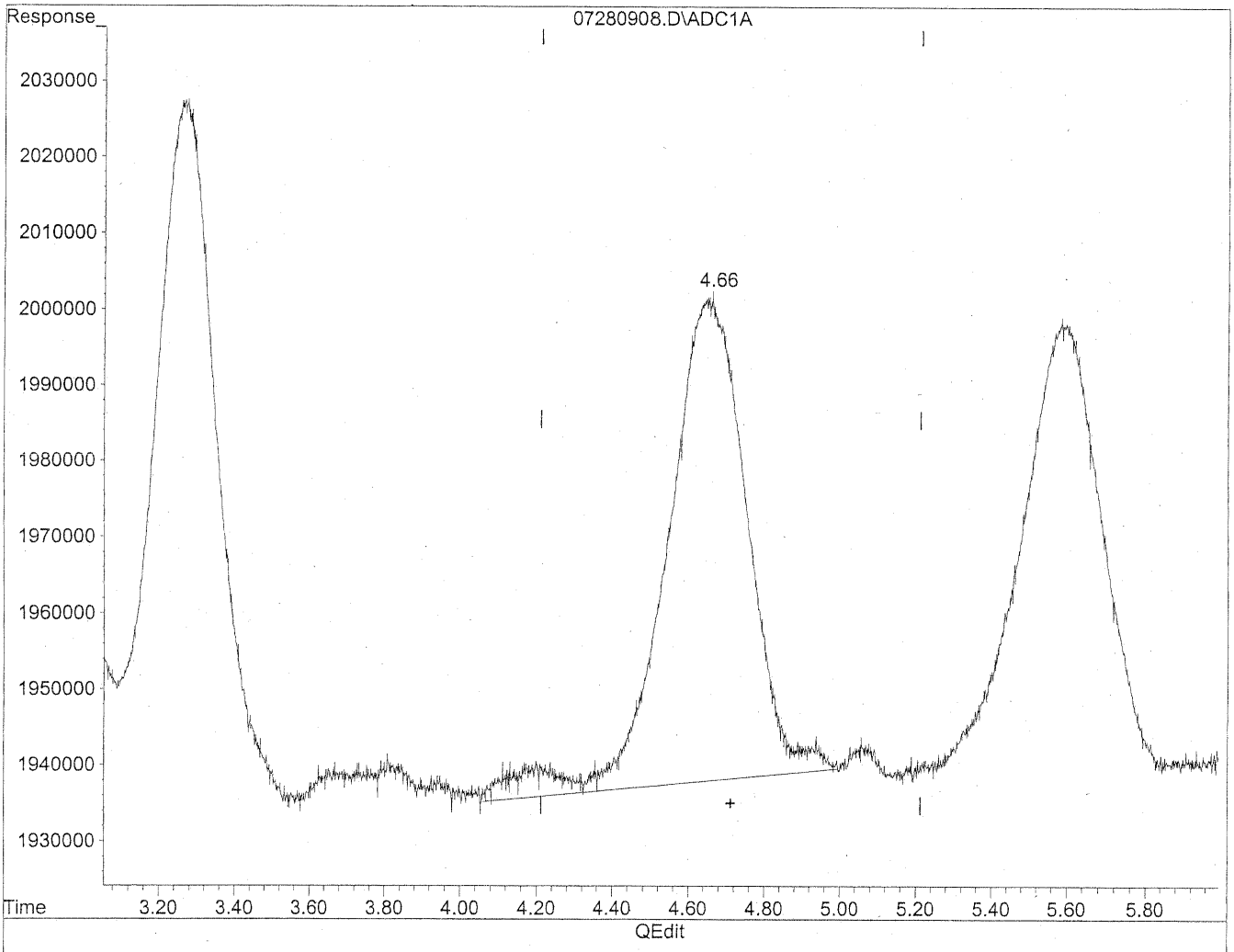
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.26	18400032	104.788	ng/ml
2)	Acetaldehyde	1.79	13737532	101.835	ng/ml
3)	Propionaldehyde	3.27	10633406	103.442	ng/ml
4)	Crotonaldehyde	4.66	9424529	85.247	ng/mlm
5)	Butyraldehyde	5.59	8463028	105.163	ng/ml
6)	Benzaldehyde	7.04	6735919	106.795	ng/mlm
7)	Isovaleraldehyde	7.89	8025579	90.529	ng/ml
8)	Valeraldehyde	8.35	7906862	95.155	ng/ml
9)	o-Tolualdehyde	9.16	5642221	104.737	ng/ml
10)	m,p-Tolualdehyde	9.48	11177259	207.507	ng/ml
11)	Hexaldehyde	11.13	6920120	103.072	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.53	4707951	90.653	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280908.D Vial: 8
Acq On : 28 Jul 2009 10:24 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:33 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.65min 85.241ng/ml

response 9423805

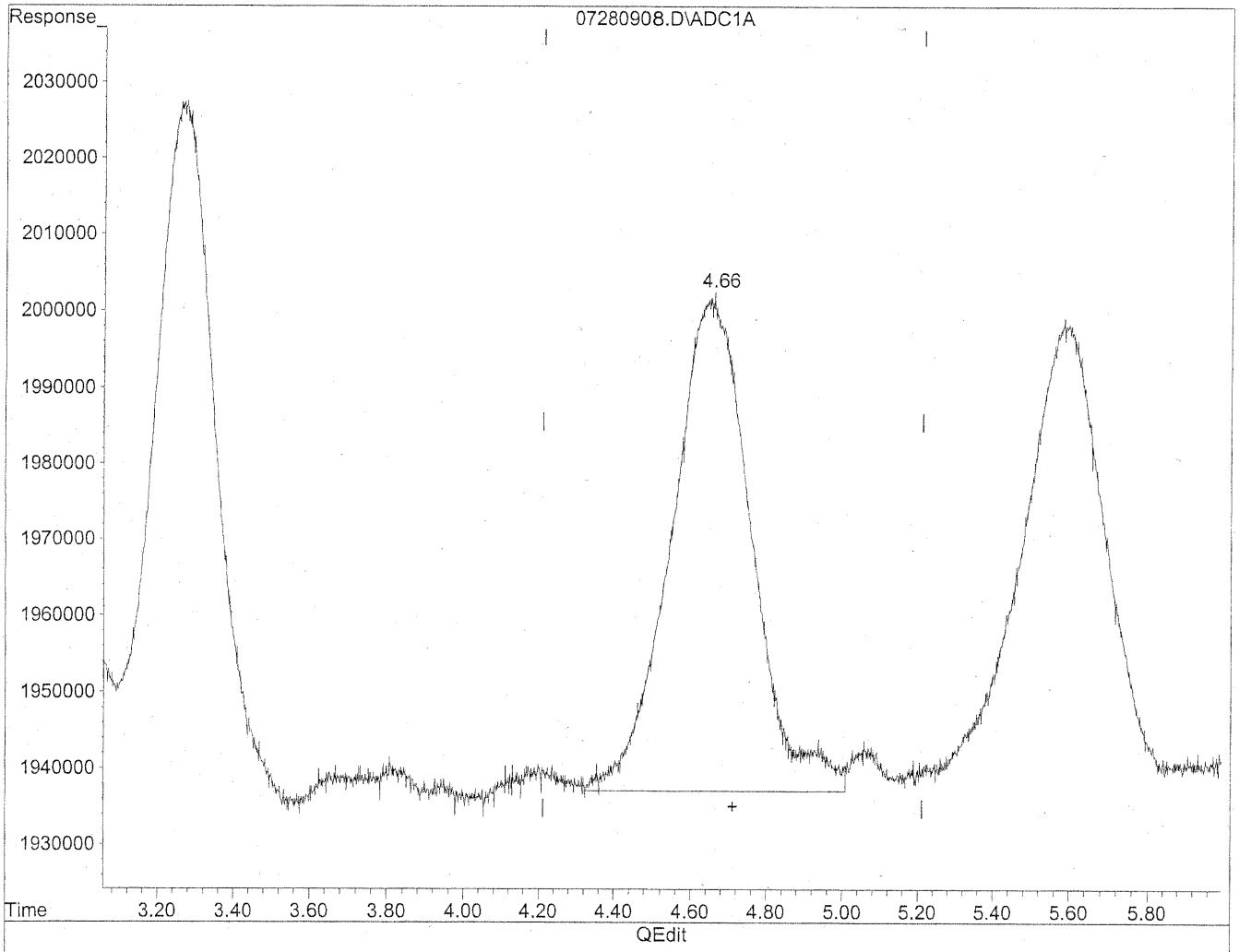
439

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280908.D Vial: 8
Acq On : 28 Jul 2009 10:24 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:33 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.66min 85.247ng/ml m

response 9424529

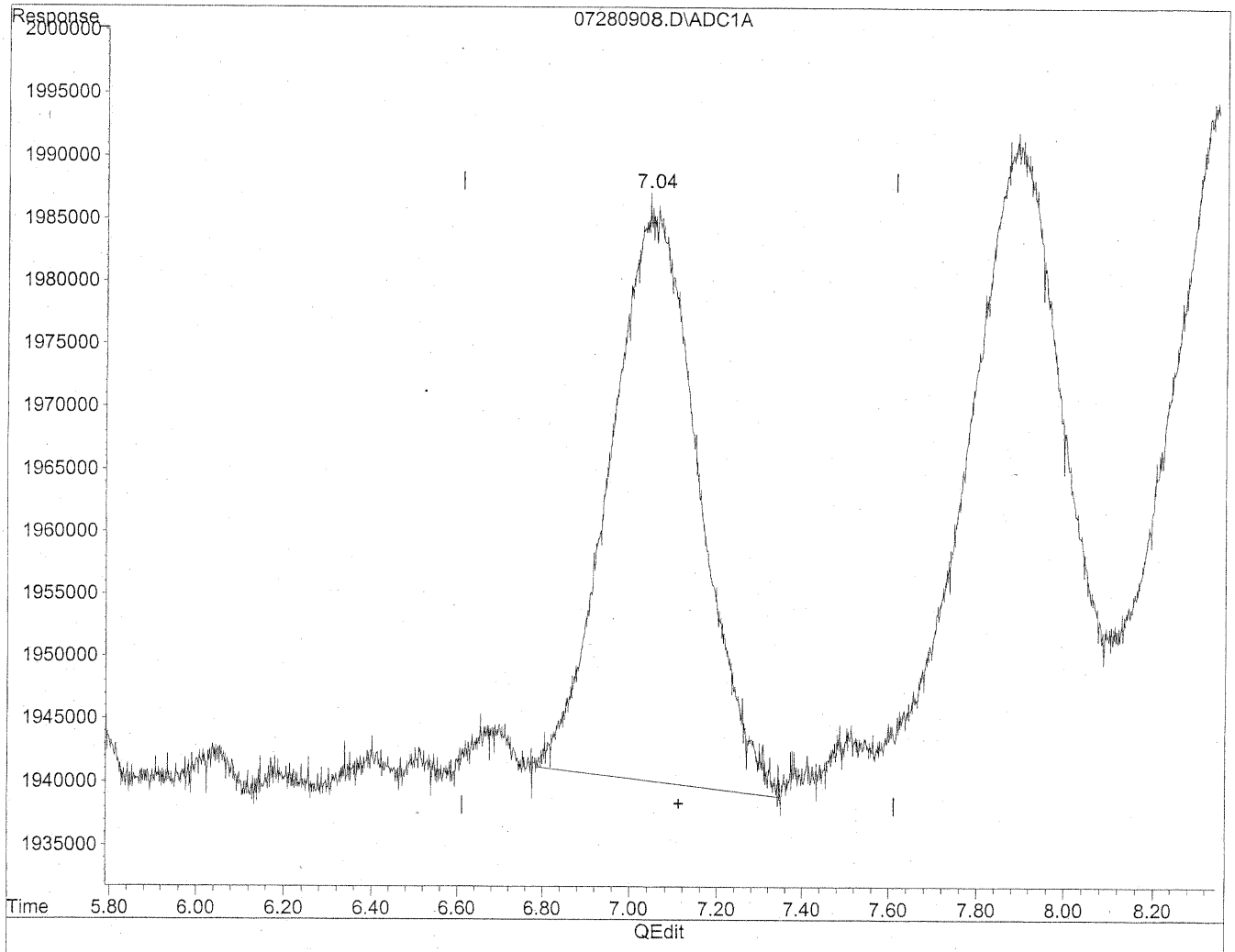
HC 7/28/09
SH
KF 7/29/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280908.D Vial: 8
Acq On : 28 Jul 2009 10:24 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:33 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

7.05min 101.515ng/ml

response 6402857

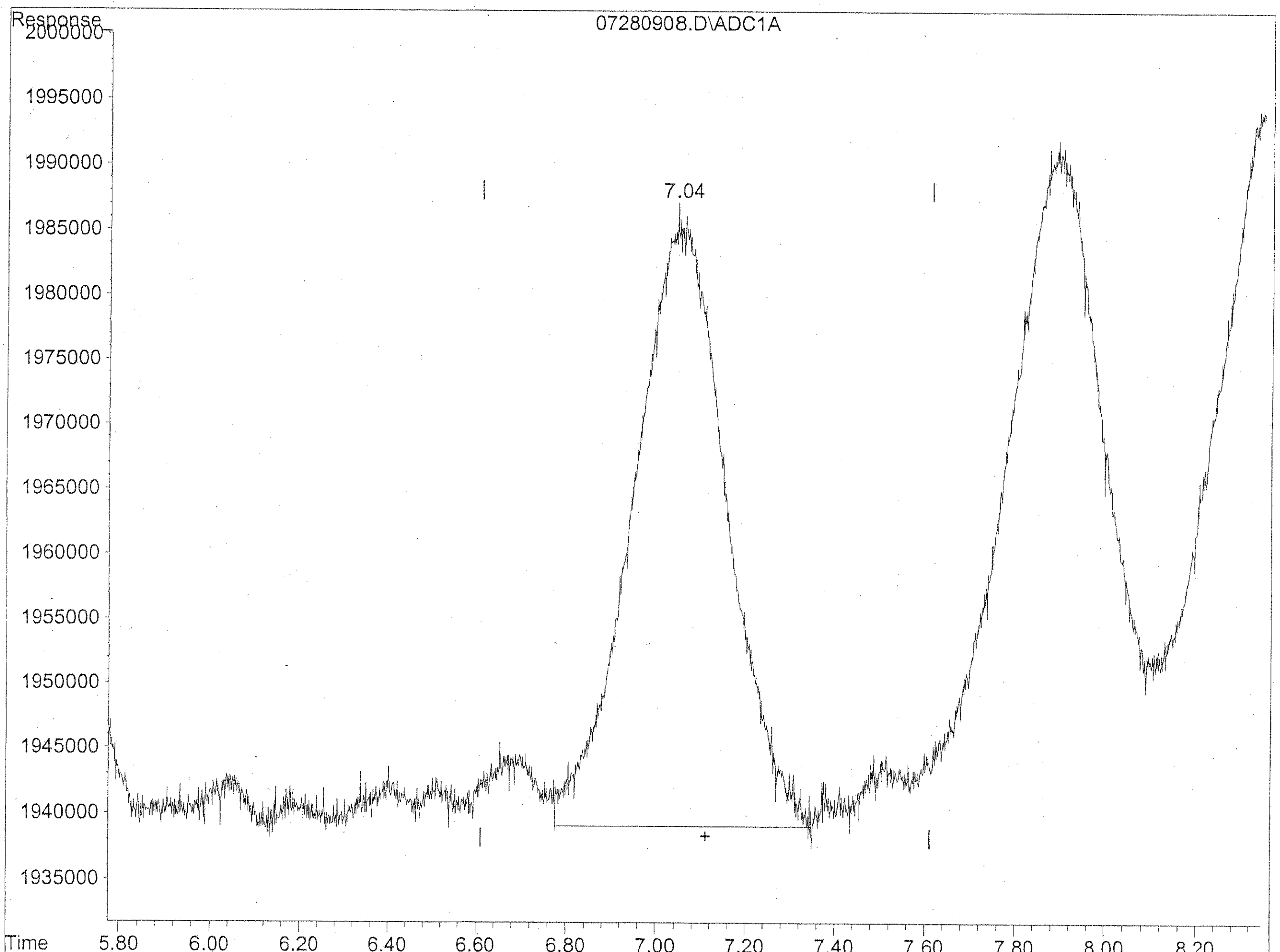
(+) = Expected Retention Time

07280908.D TO110709.M Tue Jul 28 14:34:27 2009

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2009_07\28\07280908.D        Vial: 8
Acq On    : 28 Jul 2009  10:24 am                       Operator: HC
Sample    : 100ng/ml TO11A Std S21-07270905             Inst    : LC 01
Misc      :                                             Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Jul 28 14:34 19109  Quant Results File: TO110709.RES

Method          : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title           : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update     : Tue Jul 28 15:29:52 2009
Response via    : Multiple Level Calibration
```



QEdit

(6) Benzaldehyde

7.04min 106.795ng/ml m

response 6735919

HC 7/28/09 BC

KK 7/29/09

(+) = Expected Retention Time

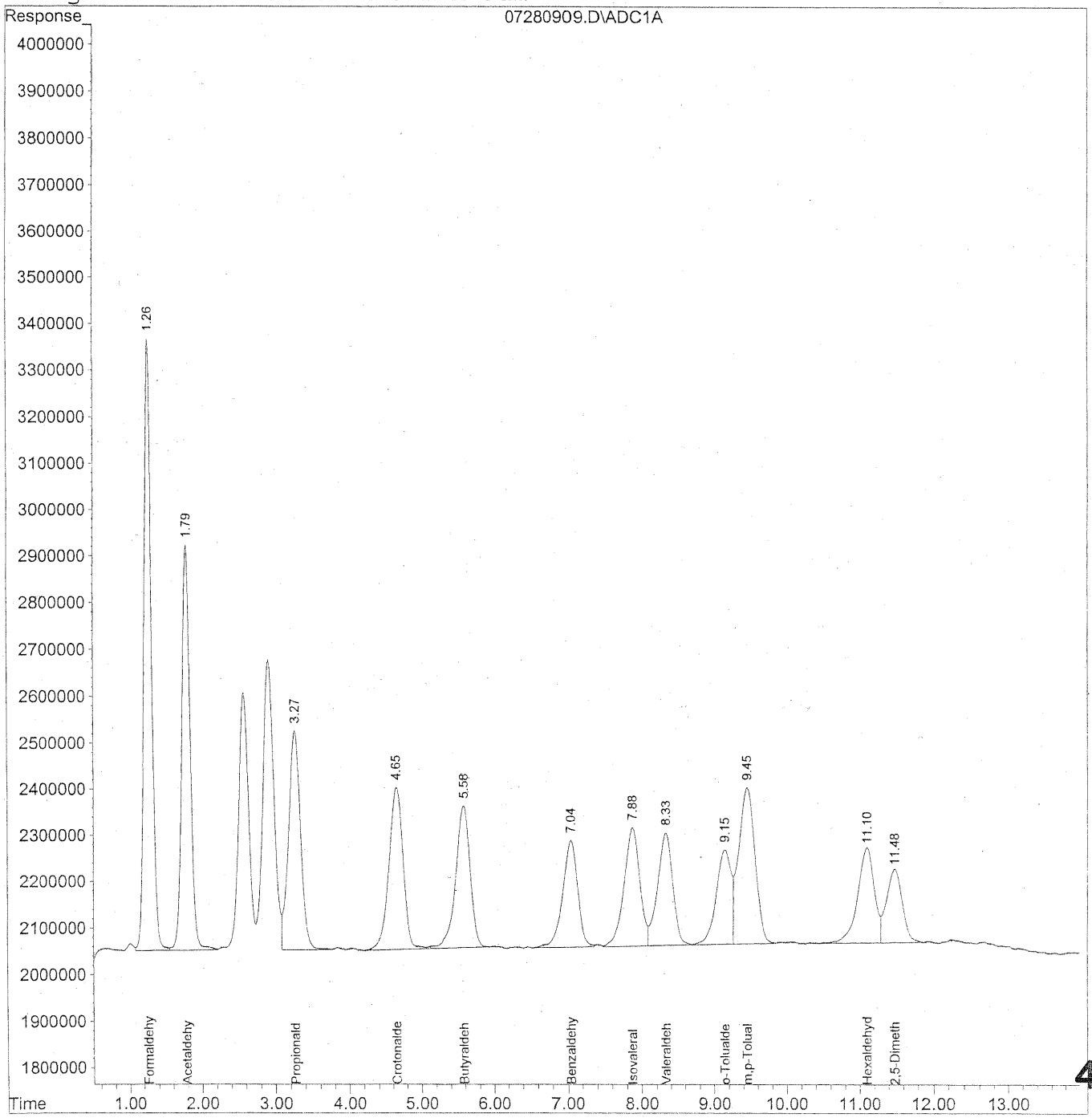
07280908.D   TO110709.M   Tue Jul 28 17:38:30 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280909.D Vial: 9
Acq On : 28 Jul 2009 10:39 am Operator: HC
Sample : 500ng/ml TO11A Std S21-07270904 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:40 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



443

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280909.D Vial: 9
Acq On : 28 Jul 2009 10:39 am Operator: HC
Sample : 500ng/ml TO11A Std S21-07270904 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:40 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.26	91593554	521.625	ng/ml
2)	Acetaldehyde	1.79	70468869	522.381	ng/ml
3)	Propionaldehyde	3.27	53468174	520.142	ng/ml
4)	Crotonaldehyde	4.65	47866960	432.968	ng/ml
5)	Butyraldehyde	5.58	43271557	537.700	ng/ml
6)	Benzaldehyde	7.04	32616313	517.119	ng/ml
7)	Isovaleraldehyde	7.88	37944016	428.013	ng/ml
8)	Valeraldehyde	8.33	35574509	428.119	ng/ml
9)	o-Tolualdehyde	9.15	29317615	544.227	ng/ml
10)	m,p-Tolualdehyde	9.46	53274975	989.053	ng/ml
11)	Hexaldehyde	11.10	32888440	489.859	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.49	23823948	458.738	ng/ml

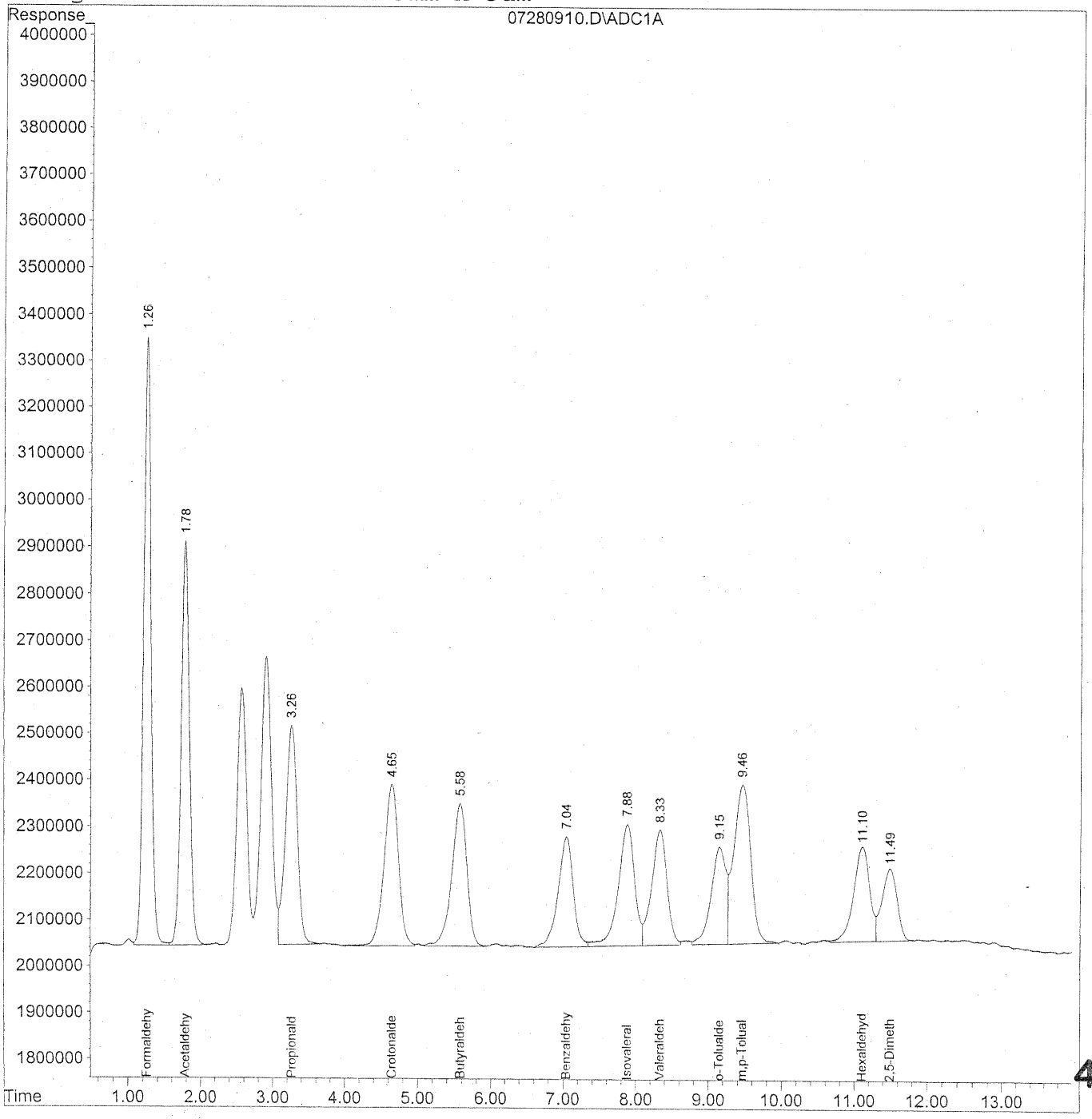
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280910.D Vial: 10
Acq On : 28 Jul 2009 10:54 am Operator: HC
Sample : 500ng/ml TO11A Std S21-07270904 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:41 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



445

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280910.D Vial: 10
Acq On : 28 Jul 2009 10:54 am Operator: HC
Sample : 500ng/ml TO11A Std S21-07270904 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:41 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.26	90711575	516.602	ng/ml
2)	Acetaldehyde	1.78	69140255	512.533	ng/ml
3)	Propionaldehyde	3.26	52850412	514.132	ng/ml
4)	Crotonaldehyde	4.65	47584179	430.411	ng/ml
5)	Butyraldehyde	5.58	43677338	542.743	ng/ml
6)	Benzaldehyde	7.04	34085310	540.409	ng/ml
7)	Isovaleraldehyde	7.88	40968120	462.125	ng/ml
8)	Valeraldehyde	8.33	36648075	441.039	ng/ml
9)	o-Tolualdehyde	9.15	29793454	553.060	ng/ml
10)	m,p-Tolualdehyde	9.46	54514161	1012.059	ng/ml
11)	Hexaldehyde	11.11	31855201	474.470	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.49	22510750	433.452	ng/ml

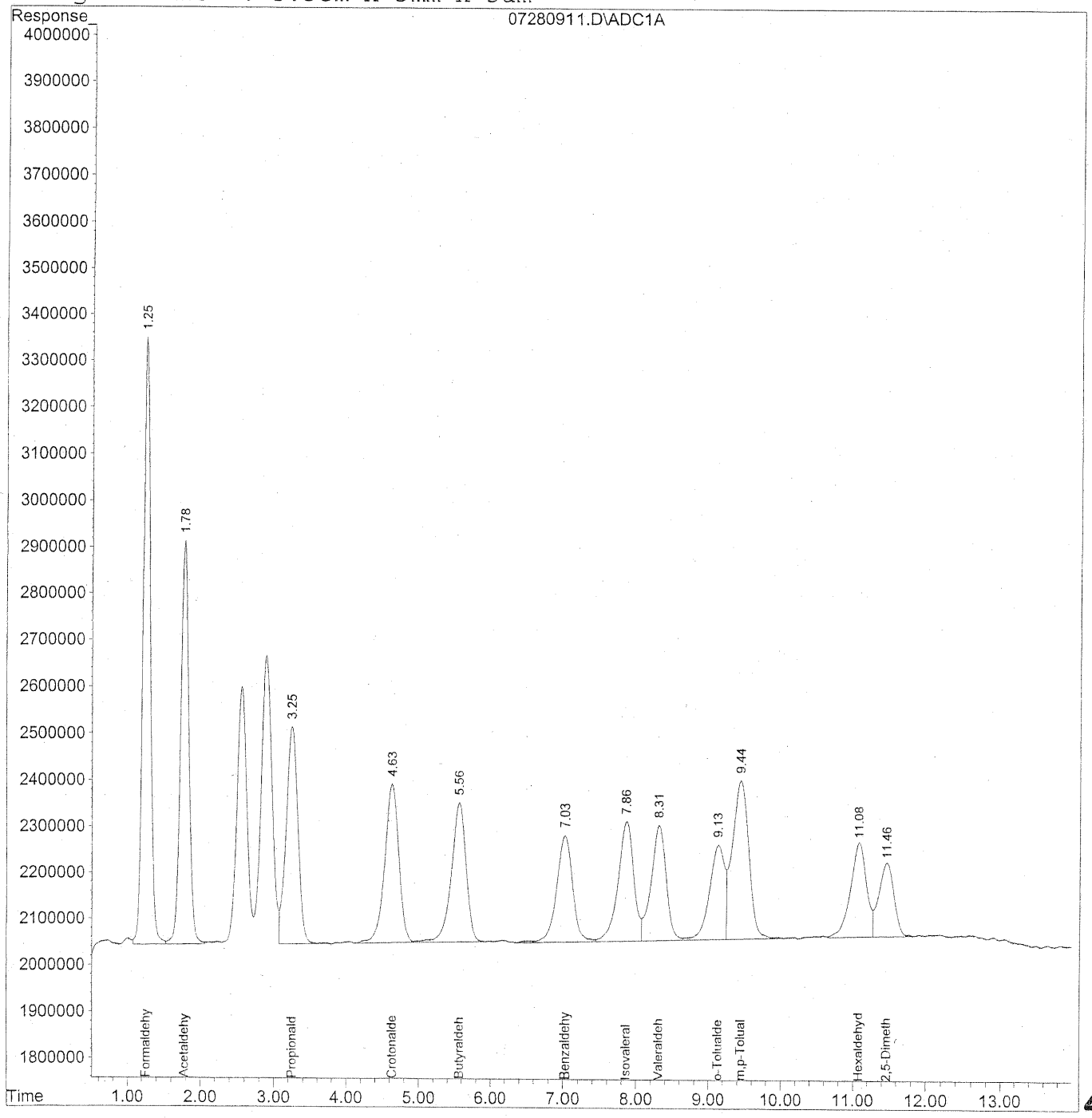
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280911.D Vial: 11
Acq On : 28 Jul 2009 11:09 am Operator: HC
Sample : 500ng/ml TO11A Std S21-07270904 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:41 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



447

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280911.D Vial: 11
Acq On : 28 Jul 2009 11:09 am Operator: HC
Sample : 500ng/ml TO11A Std S21-07270904 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:41 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

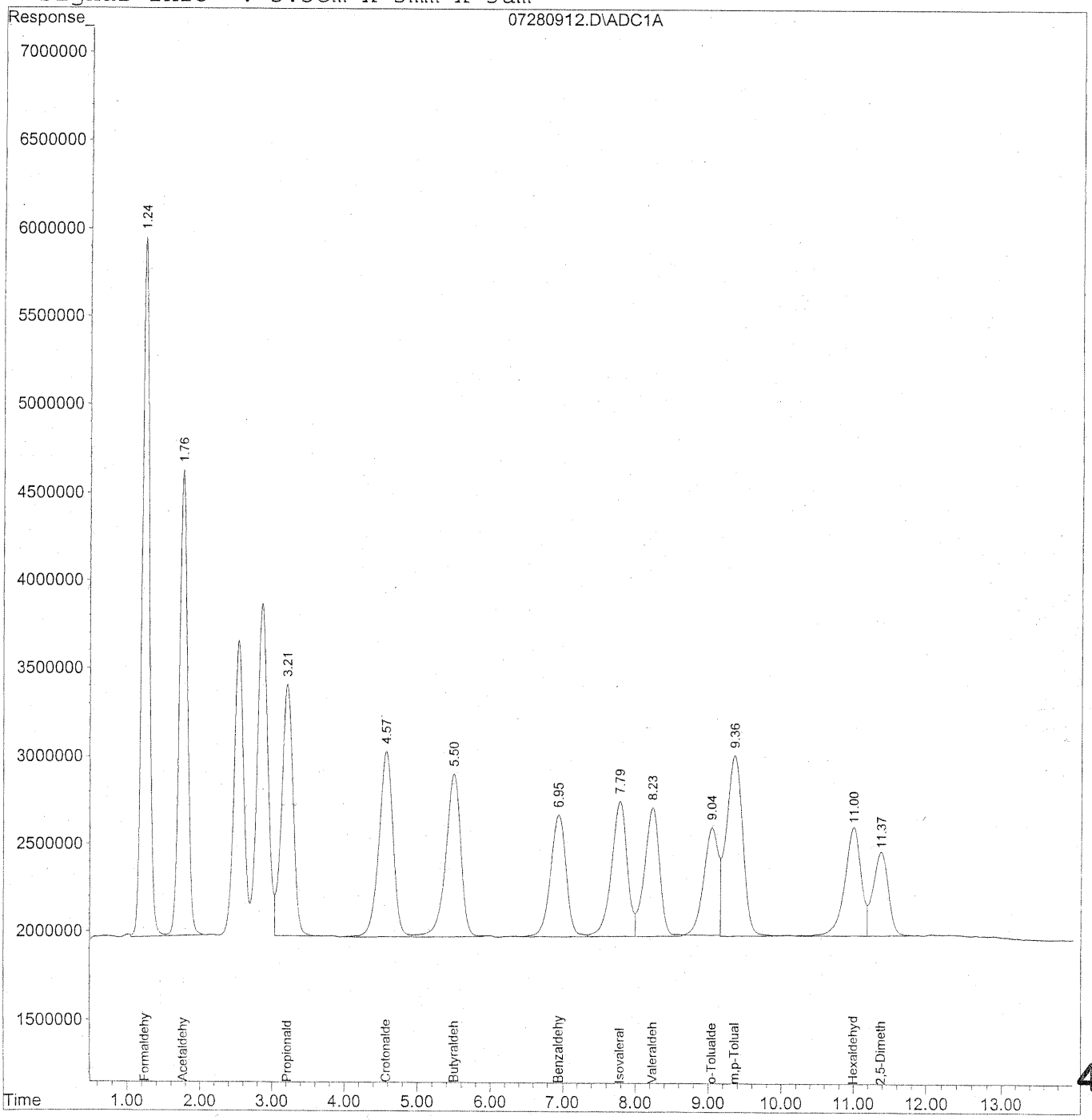
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.25	91399555	520.520	ng/ml
2)	Acetaldehyde	1.78	69908753	518.229	ng/ml
3)	Propionaldehyde	3.25	52190620	507.713	ng/ml
4)	Crotonaldehyde	4.63	46362546	419.361	ng/ml
5)	Butyraldehyde	5.56	43673214	542.691	ng/ml
6)	Benzaldehyde	7.03	34084716	540.400	ng/ml
7)	Isovaleraldehyde	7.87	39175205	441.901	ng/ml
8)	Valeraldehyde	8.31	36501988	439.281	ng/ml
9)	o-Tolualdehyde	9.13	30169058	560.032	ng/ml
10)	m,p-Tolualdehyde	9.44	54668231	1014.919	ng/ml
11)	Hexaldehyde	11.08	32179520	479.300	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.46	23309464	448.831	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Data File : J:\LC01\DATA\TO11\2009_07\28\07280912.D Vial: 12
Acq On : 28 Jul 2009 11:24 am Operator:
Sample : 1500ng/ml TO11A Std S21-07270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:45 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



449

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280912.D Vial: 12
Acq On : 28 Jul 2009 11:24 am Operator:
Sample : 1500ng/ml TO11A Std S21-07270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:45 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

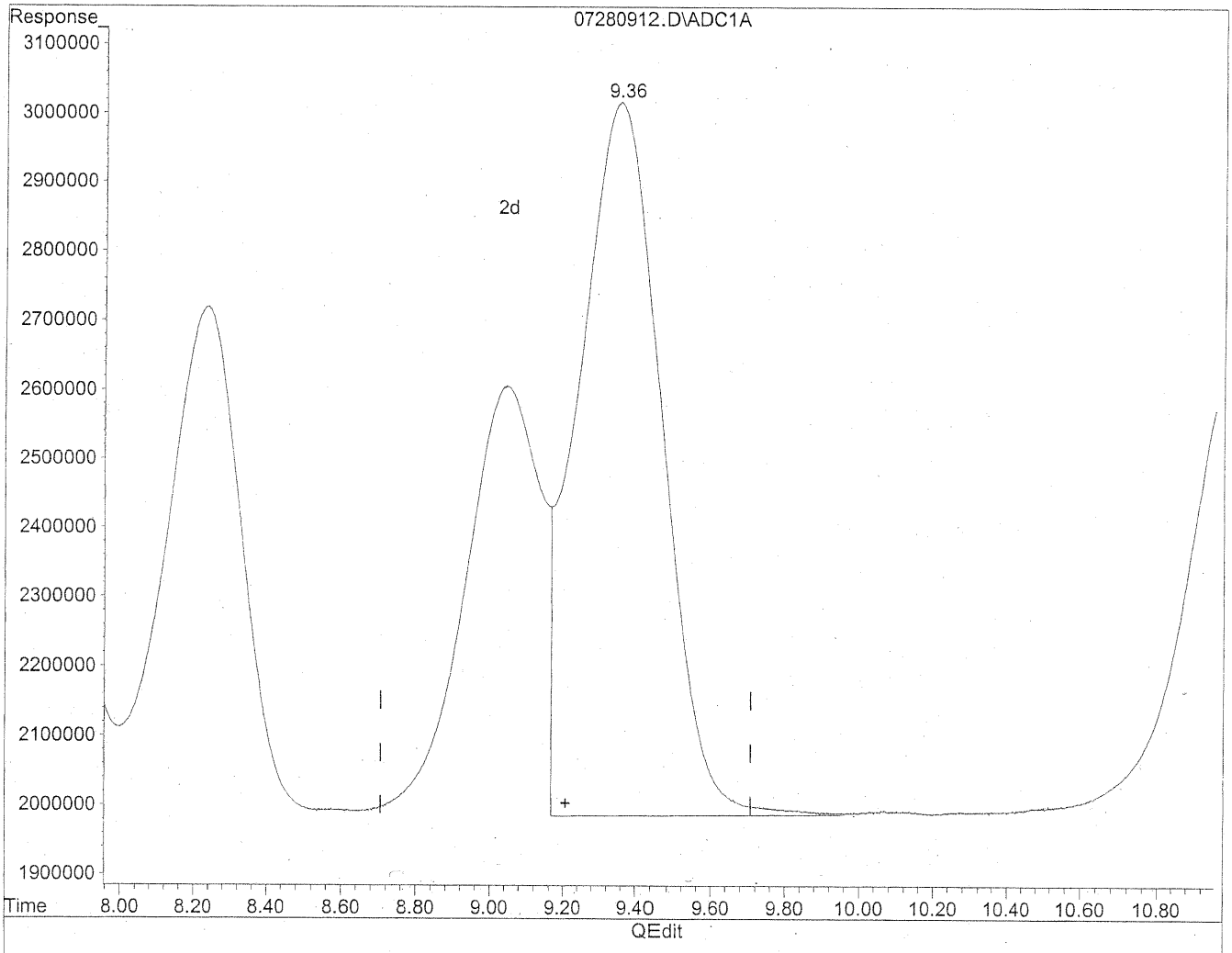
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.24	275380897	1568.292	ng/ml
2)	Acetaldehyde	1.76	209374751	1552.082	ng/ml
3)	Propionaldehyde	3.21	159030091	1547.054	ng/ml
4)	Crotonaldehyde	4.57	143227783	1295.530	ng/ml
5)	Butyraldehyde	5.50	134132687	1666.757	ng/ml
6)	Benzaldehyde	6.95	98878868	1567.685	ng/ml
7)	Isovaleraldehyde	7.78	115866442	1306.987	ng/ml
8)	Valeraldehyde	8.23	107104204	1288.938	ng/ml
9)	o-Tolualdehyde	9.05	86339652	1602.734	ng/mlm
10)	m,p-Tolualdehyde	9.35	162946532	3025.113	ng/ml
11)	Hexaldehyde	11.00f	98895406	1473.005	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.37	69932636	1346.576	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280912.D Vial: 12
Acq On : 28 Jul 2009 11:24 am Operator: HC
Sample : 1500ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:41 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.35min 3024.797ng/ml

response 162946532

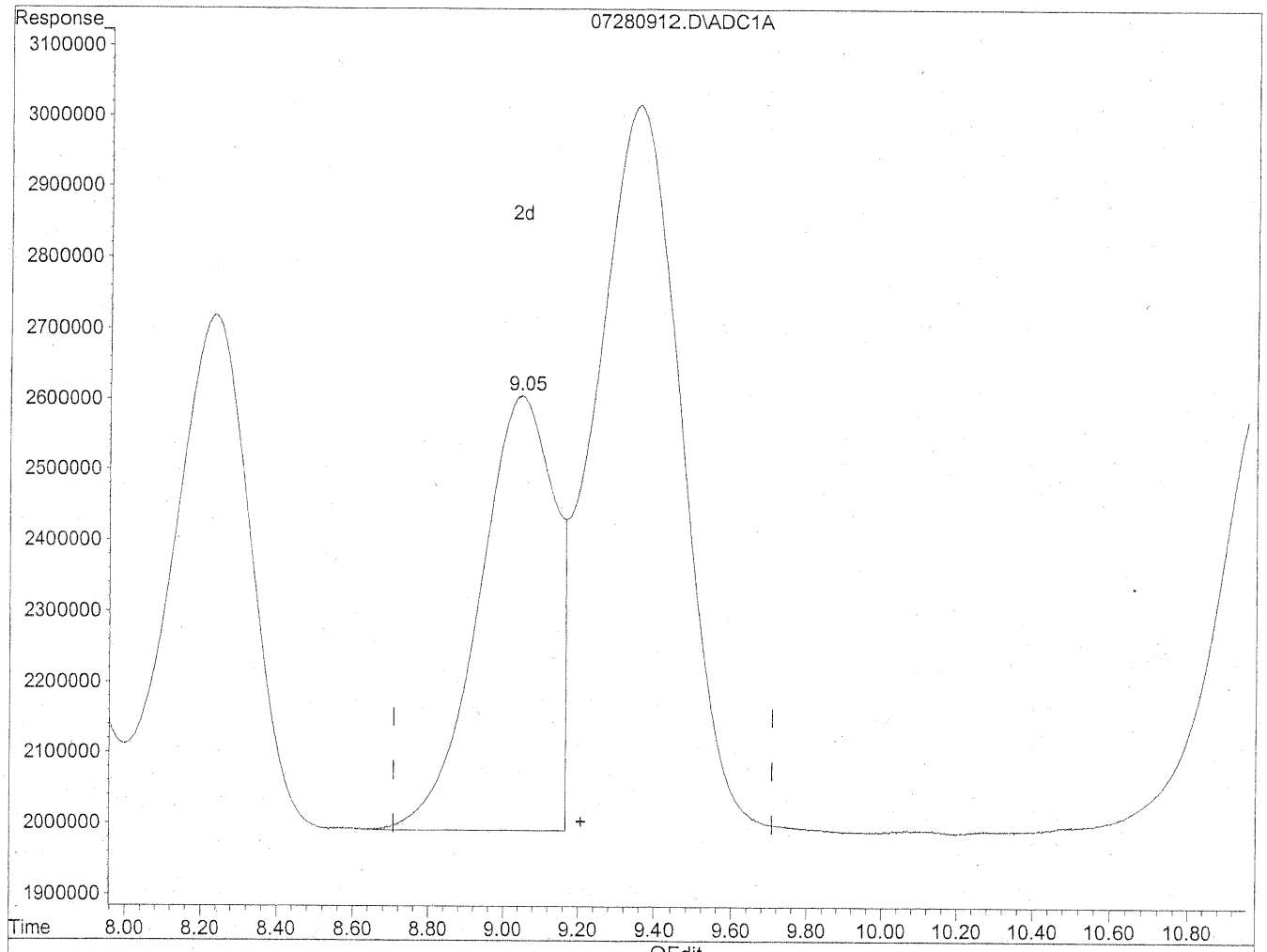
451

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280912.D Vial: 12
Acq On : 28 Jul 2009 11:24 am Operator: HC
Sample : 1500ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:41 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.05min 1602.734ng/ml m

response 86339652

HC 7/28/09 MP

7/29/09

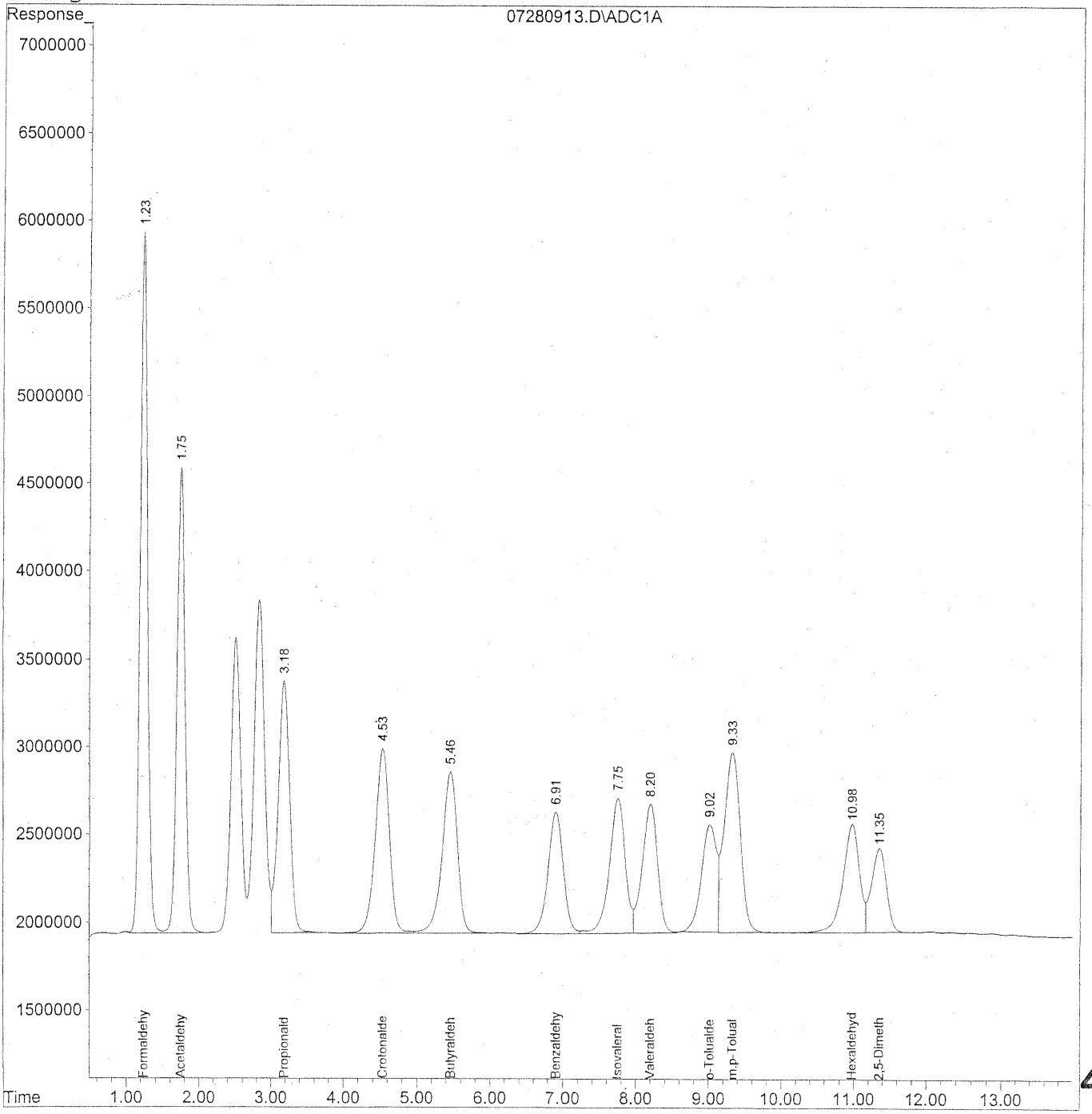
452

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_07\28\07280913.D Vial: 13
Acq On : 28 Jul 2009 11:39 am Operator: HC
Sample : 1500ng/ml TO11A Std S21-07270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:47 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



453

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280913.D Vial: 13
Acq On : 28 Jul 2009 11:39 am Operator: HC
Sample : 1500ng/ml TO11A Std S21-07270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:47 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

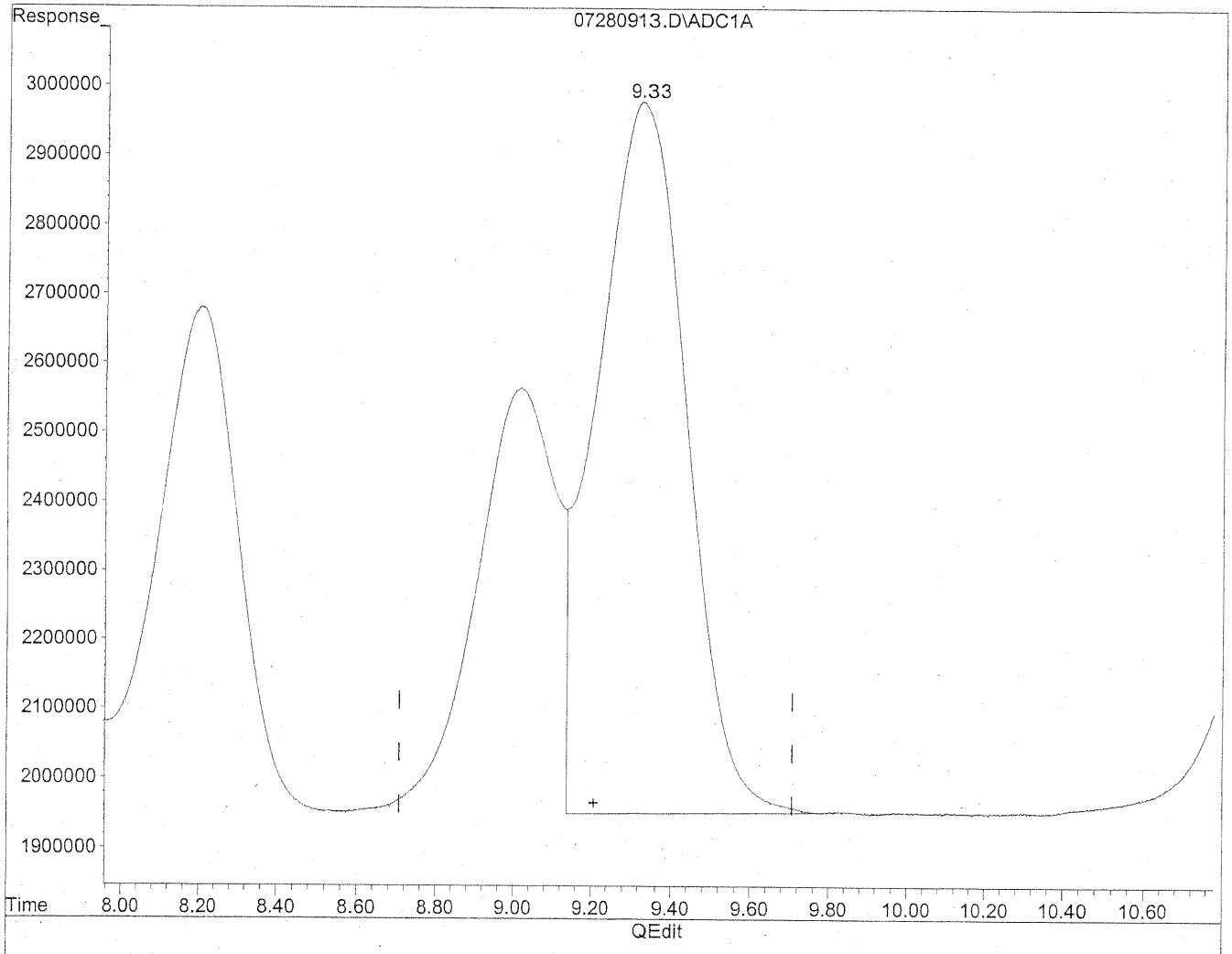
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.23	274724982	1564.557	ng/ml
2)	Acetaldehyde	1.75	209301649	1551.540	ng/ml
3)	Propionaldehyde	3.18	158919579	1545.979	ng/ml
4)	Crotonaldehyde	4.53	142112419	1285.442	ng/ml
5)	Butyraldehyde	5.46	132549734	1647.087	ng/ml
6)	Benzaldehyde	6.91	98183657	1556.663	ng/ml
7)	Isovaleraldehyde	7.75	116723586	1316.656	ng/ml
8)	Valeraldehyde	8.20	107107592	1288.979	ng/ml
9)	o-Tolualdehyde	9.02	85940120	1595.318	ng/mlm
10)	m,p-Tolualdehyde	9.33	161094009	2990.721	ng/ml
11)	Hexaldehyde	10.98f	98090122	1461.011	ng/mlm
12)	2,5-Dimethylbenzaldehyde	11.35	68873541	1326.183	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280913.D Vial: 13
Acq On : 28 Jul 2009 11:39 am Operator: HC
Sample : 1500ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:46 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.33min 2990.409ng/ml

response 161094009

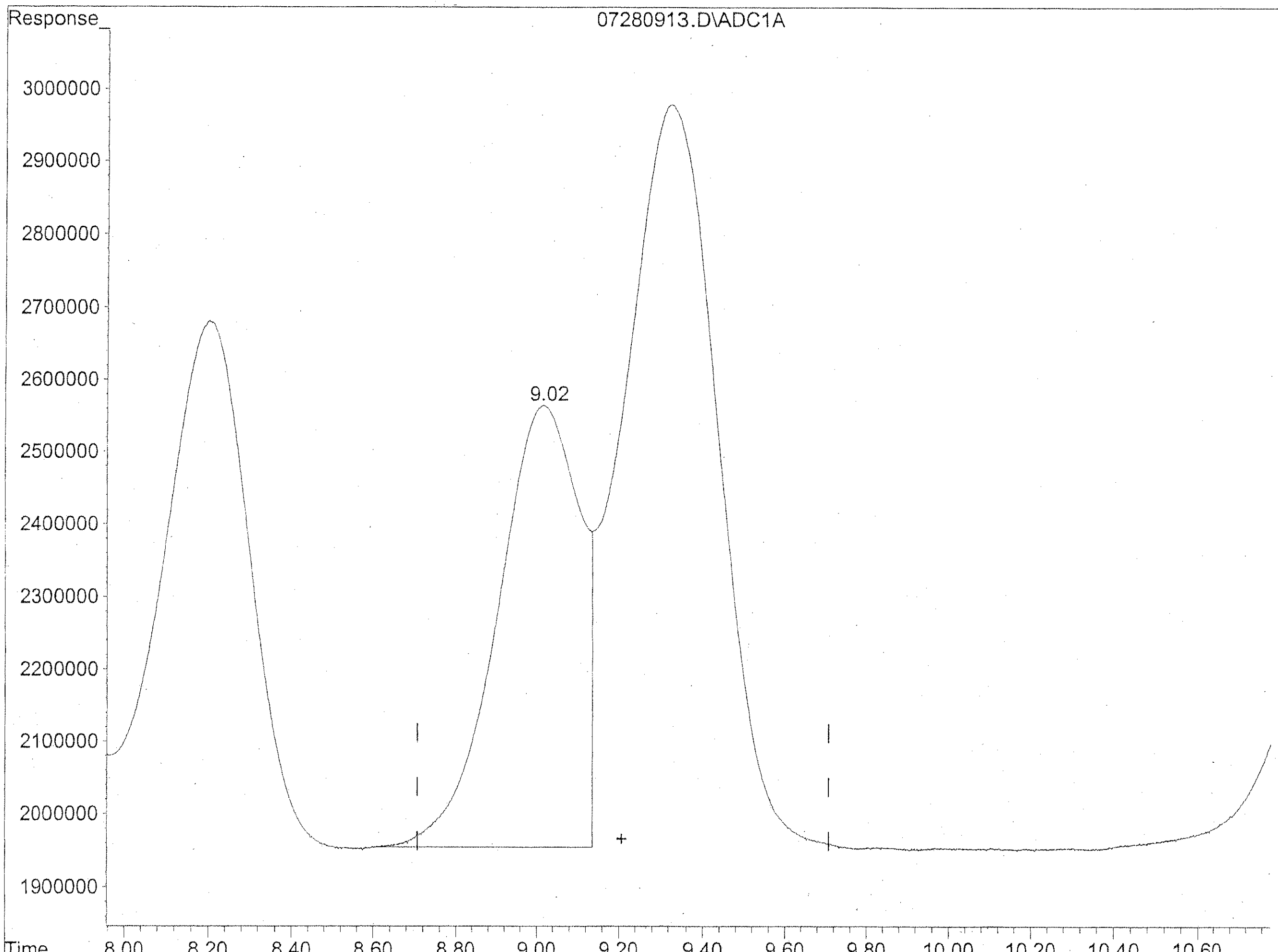
455

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280913.D Vial: 13
Acq On : 28 Jul 2009 11:39 am Operator: HC
Sample : 1500ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:46 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



QEdit

(9) o-Tolualdehyde

9.02min 1595.318ng/ml m

response 85940120

HC 7/28/09 MP

KR 7/29/09

456

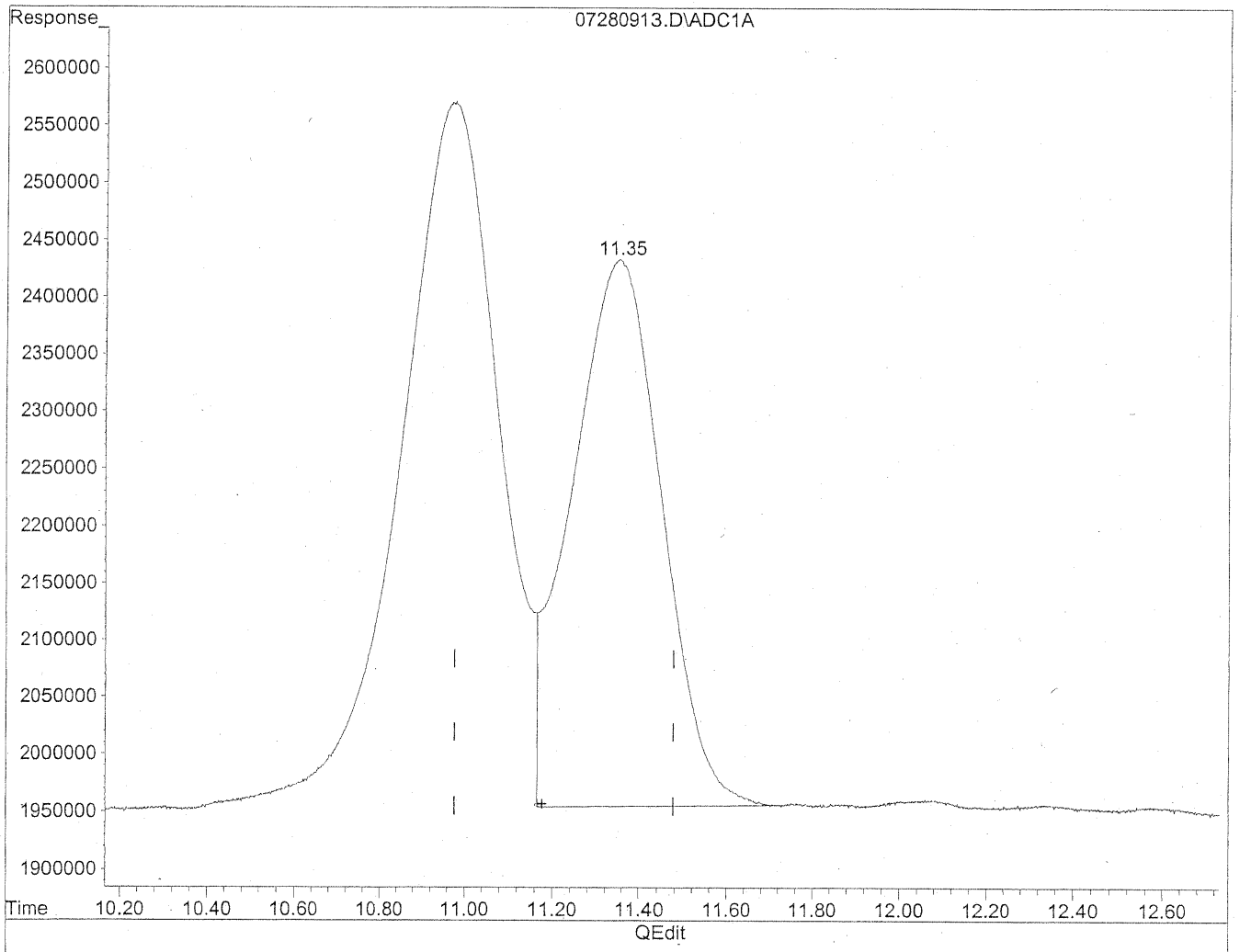
(+) = Expected Retention Time

07280913.D TO110709.M Tue Jul 28 14:47:19 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280913.D Vial: 13
Acq On : 28 Jul 2009 11:39 am Operator: HC
Sample : 1500ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:46 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

11.35min 1025.842ng/ml

response 68873541

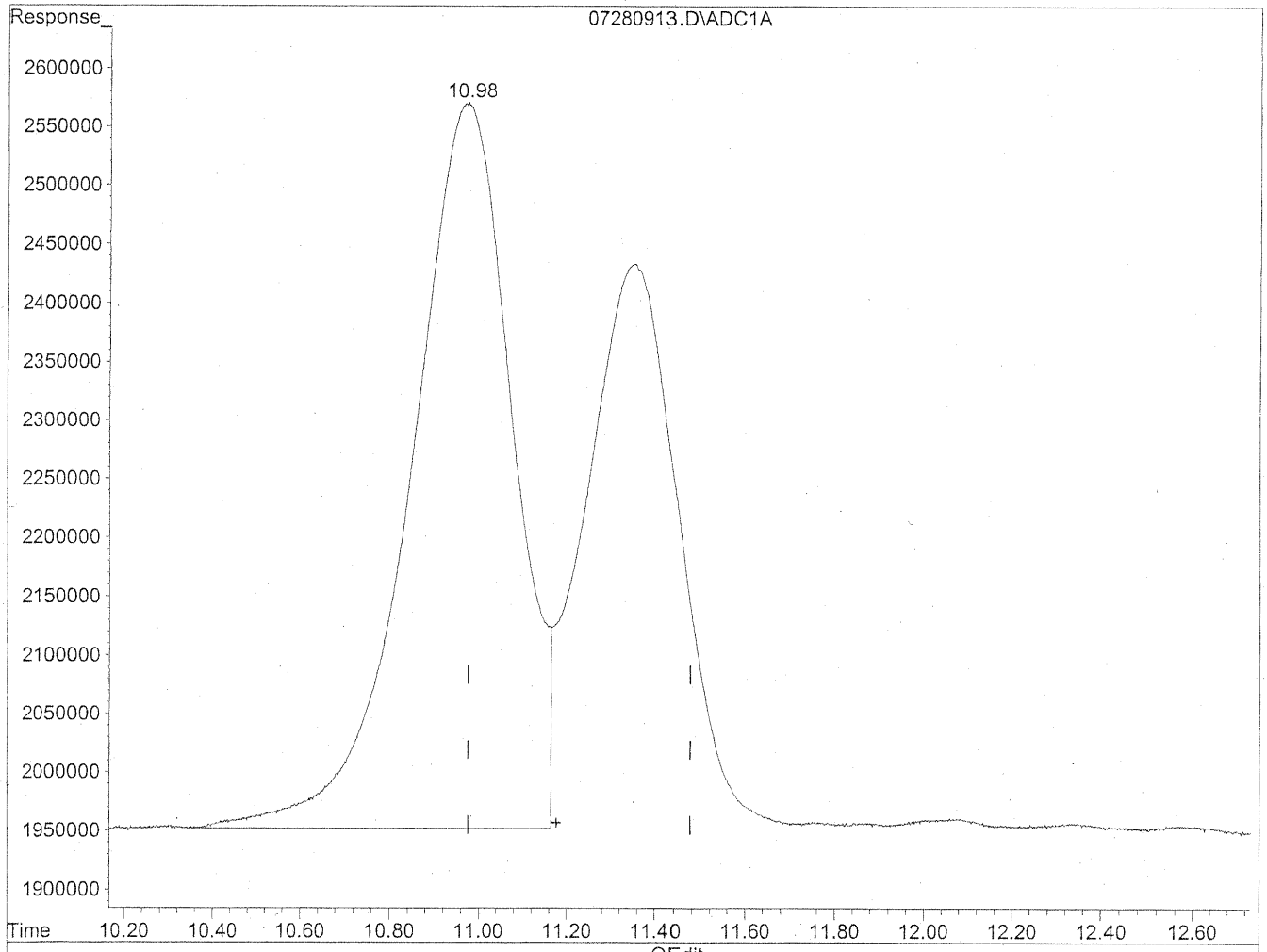
457

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280913.D Vial: 13
Acq On : 28 Jul 2009 11:39 am Operator: HC
Sample : 1500ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:46 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



QEdit

(11) Hexaldehyde

10.98min 1461.011ng/ml m

response 98090122

458

(+) = Expected Retention Time

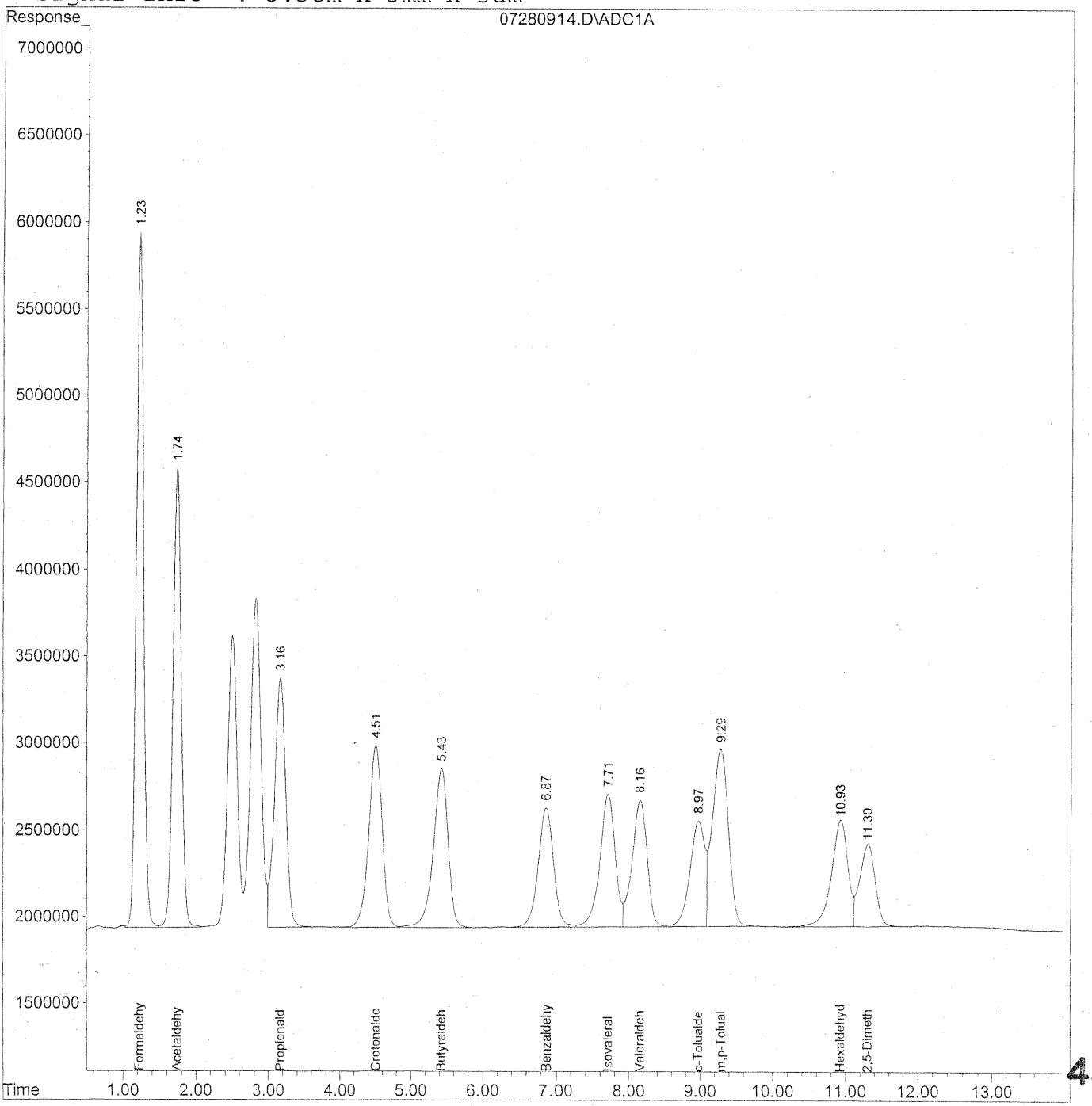
07280913.D TO110709.M Tue Jul 28 14:47:35 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280914.D Vial: 14
Acq On : 28 Jul 2009 11:54 am Operator: HC
Sample : 1500ng/ml TO11A Std S21-07270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:48 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



459

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280914.D Vial: 14
Acq On : 28 Jul 2009 11:54 am Operator: HC
Sample : 1500ng/ml TO11A Std S21-07270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:48 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.23	273895978	1526.977	ng/ml
2)	Acetaldehyde	1.74	208465321	1519.240	ng/ml
3)	Propionaldehyde	3.16	158125683	1515.036	ng/ml
4)	Crotonaldehyde	4.50	139629551	1360.269	ng/ml
5)	Butyraldehyde	5.43	131425702	1556.792	ng/ml
6)	Benzaldehyde	6.87	97652643	1519.543	ng/ml
7)	Isovaleraldehyde	7.71	114690000	1377.928	ng/ml
8)	Valeraldehyde	8.16	105937177	1371.355	ng/ml
9)	o-Tolualdehyde	8.97	87824227	1580.089	ng/ml
10)	m,p-Tolualdehyde	9.28	159292531	2961.857	ng/ml
11)	Hexaldehyde	10.93	98846718	1491.666	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.31	70224395	1435.357	ng/ml

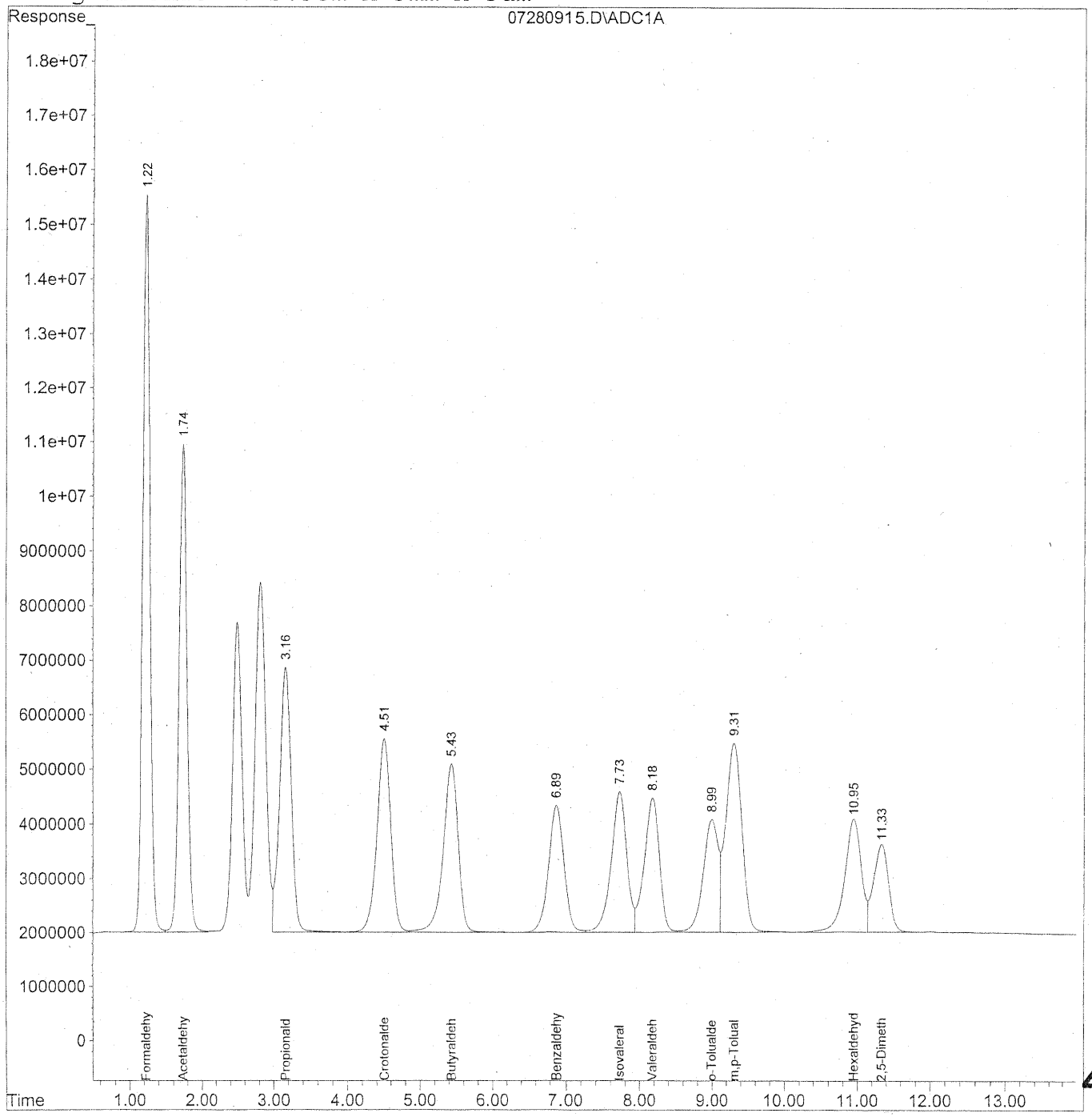
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280915.D Vial: 15
Acq On : 28 Jul 2009 12:09 pm Operator: HC
Sample : 5000ng/ml TO11A Std S21-07270902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:48 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



461

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280915.D Vial: 15
Acq On : 28 Jul 2009 12:09 pm Operator: HC
Sample : 5000ng/ml TO11A Std S21-07270902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:48 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	928364658	5175.655	ng/ml
2)	Acetaldehyde	1.74	706170560	5146.384	ng/ml
3)	Propionaldehyde	3.16	539067854	5164.924	ng/ml
4)	Crotonaldehyde	4.51	476268543	4639.802	ng/ml
5)	Butyraldehyde	5.43	446392739	5287.707	ng/ml
6)	Benzaldehyde	6.89	328286106	5108.361	ng/ml
7)	Isovaleraldehyde	7.73	388247386	4664.549	ng/ml
8)	Valeraldehyde	8.18	357832844	4632.141	ng/ml
9)	o-Tolualdehyde	8.99	298513860	5370.710	ng/ml
10)	m,p-Tolualdehyde	9.31	545640330	10145.539	ng/ml
11)	Hexaldehyde	10.95	332315493	5014.874	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.33	235692401	4817.453	ng/ml

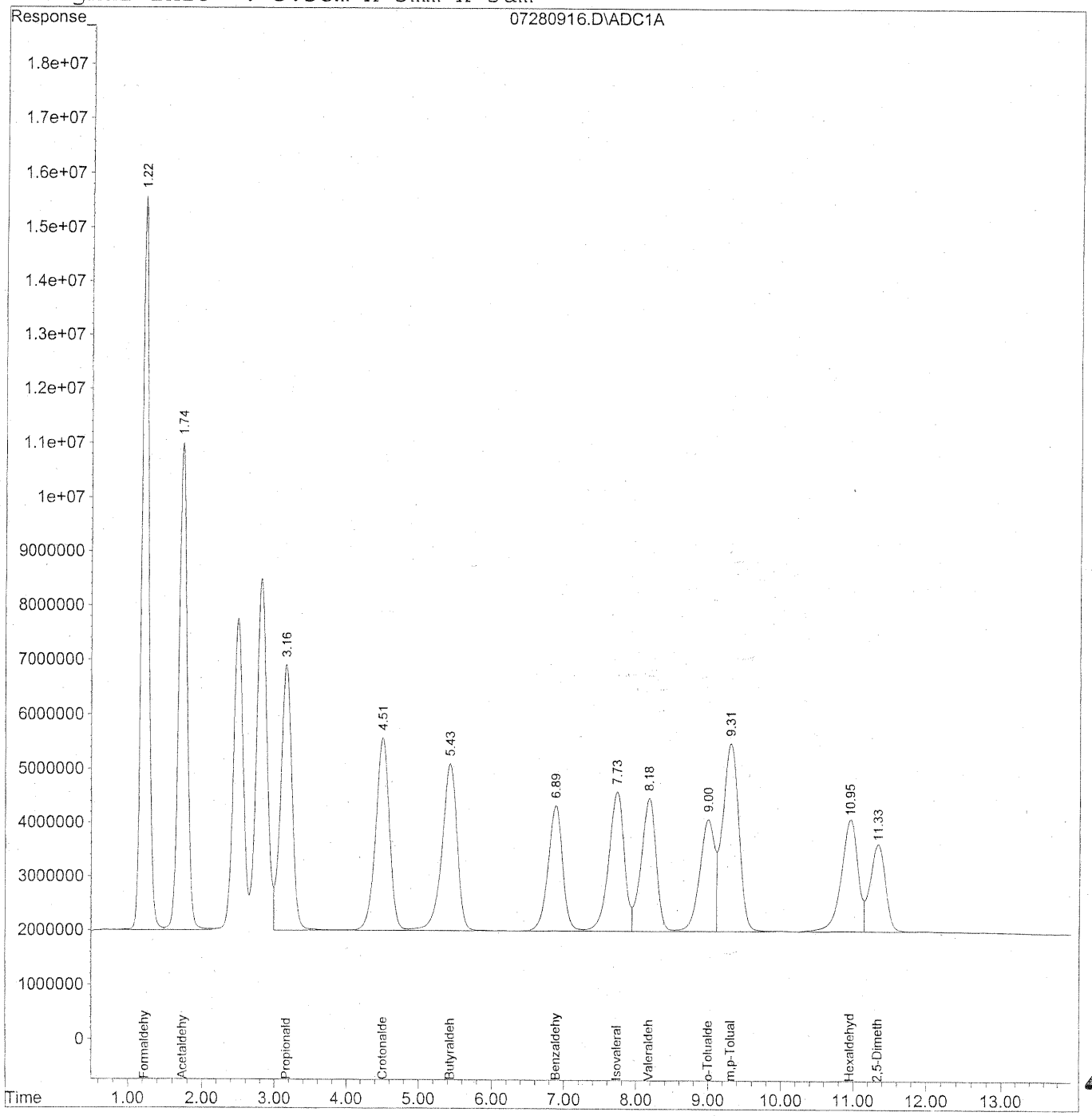
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280916.D Vial: 16
Acq On : 28 Jul 2009 12:24 pm Operator: HC
Sample : 5000ng/ml TO11A Std S21-07270902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:48 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



463

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280916.D Vial: 16
Acq On : 28 Jul 2009 12:24 pm Operator: HC
Sample : 5000ng/ml TO11A Std S21-07270902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:48 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	925768000	5161.179	ng/ml
2)	Acetaldehyde	1.74	708552415	5163.742	ng/ml
3)	Propionaldehyde	3.16	540133923	5175.139	ng/ml
4)	Crotonaldehyde	4.51	477844499	4655.155	ng/ml
5)	Butyraldehyde	5.43	446568052	5289.783	ng/ml
6)	Benzaldehyde	6.89	328413551	5110.344	ng/ml
7)	Isovaleraldehyde	7.73	388941560	4672.889	ng/ml
8)	Valeraldehyde	8.18	359676615	4656.008	ng/ml
9)	o-Tolualdehyde	9.00	300077384	5398.840	ng/ml
10)	m,p-Tolualdehyde	9.31	547211501	10174.753	ng/ml
11)	Hexaldehyde	10.95	333701808	5035.794	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.34	237108293	4846.394	ng/ml

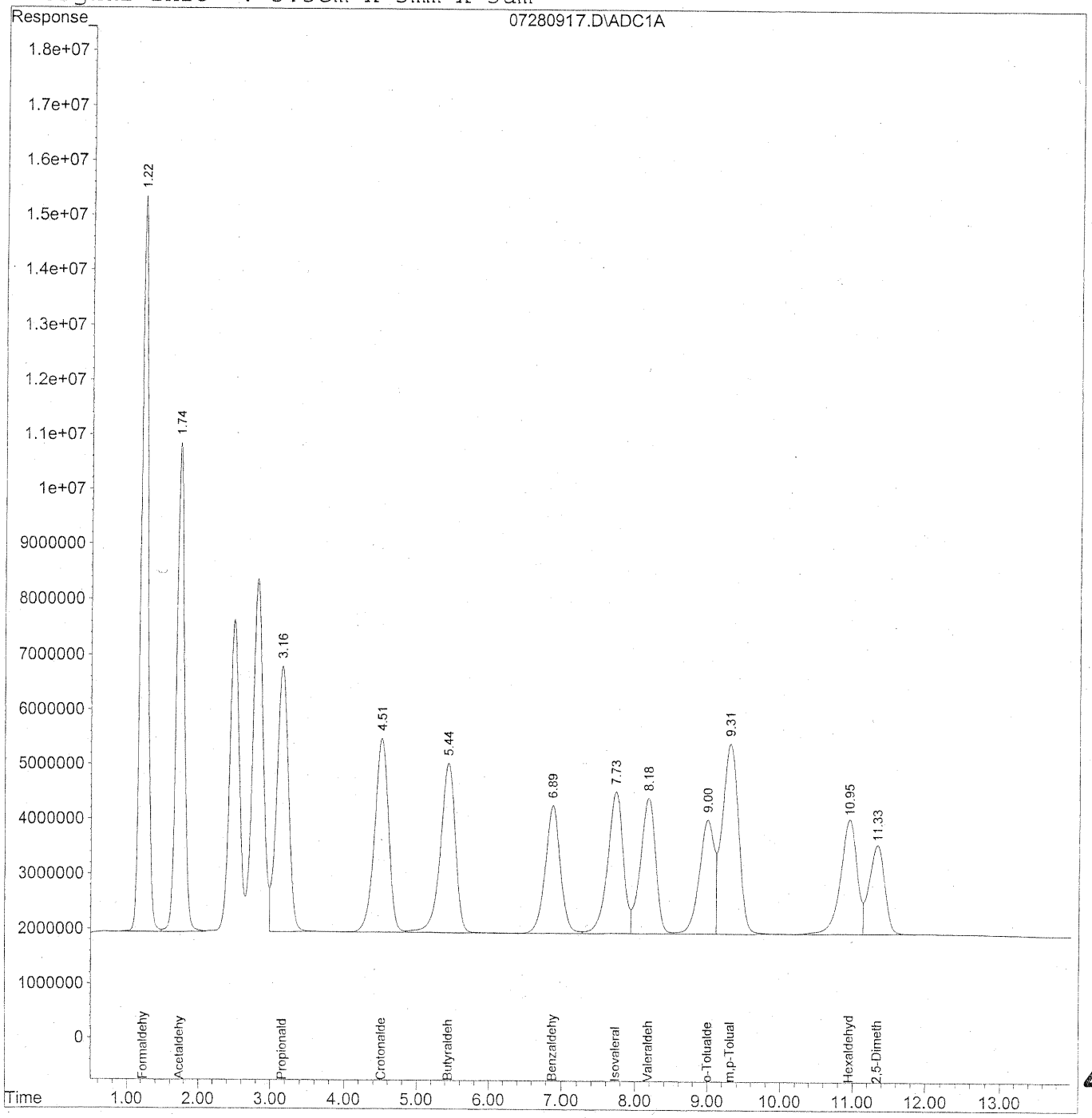
(f)=RT Delta > 1/2 Window (m)=manual int.
07280916.D TO110709.M Tue Jul 28 17:20:20 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280917.D Vial: 17
Acq On : 28 Jul 2009 12:39 pm Operator: HC
Sample : 5000ng/ml TO11A Std S21-07270902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:49 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



465

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280917.D Vial: 17
Acq On : 28 Jul 2009 12:39 pm Operator: HC
Sample : 5000ng/ml TO11A Std S21-07270902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:49 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.23	918424042	5120.236	ng/ml
2)	Acetaldehyde	1.74	702791887	5121.761	ng/ml
3)	Propionaldehyde	3.16	531675082	5094.093	ng/ml
4)	Crotonaldehyde	4.51	471954575	4597.775	ng/ml
5)	Butyraldehyde	5.44	443441833	5252.752	ng/ml
6)	Benzaldehyde	6.89	327762901	5100.219	ng/ml
7)	Isovaleraldehyde	7.73	386992833	4649.476	ng/ml
8)	Valeraldehyde	8.18	356464469	4614.427	ng/ml
9)	o-Tolualdehyde	9.00	297374461	5350.211	ng/ml
10)	m,p-Tolualdehyde	9.31	544331756	10121.207	ng/ml
11)	Hexaldehyde	10.95	332038452	5010.693	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.33	236428207	4832.493	ng/ml

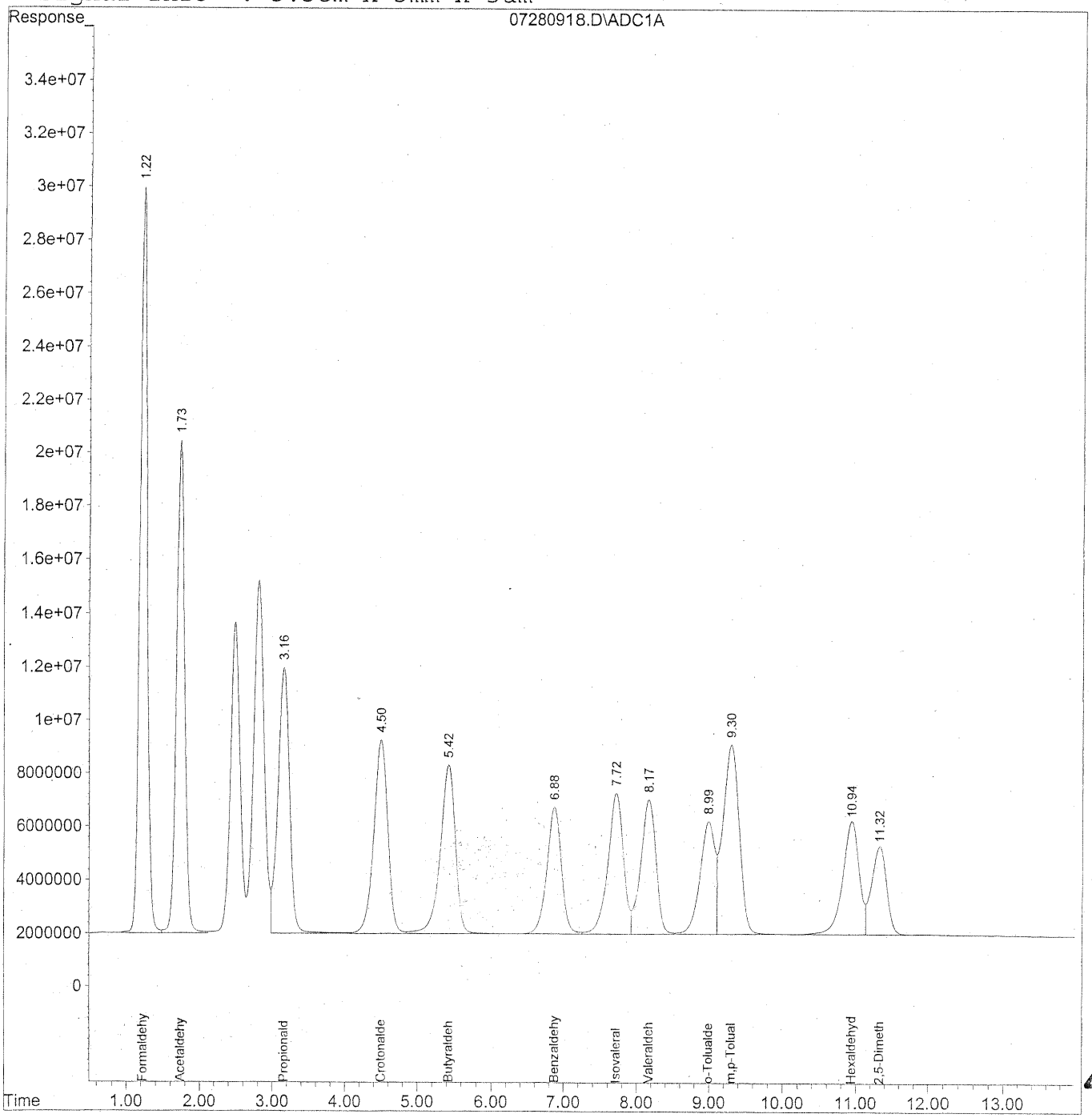
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280918.D Vial: 18
Acq On : 28 Jul 2009 12:54 pm Operator: HC
Sample : 10000ng/ml TO11A Std S21-07270901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:49 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



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Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280918.D Vial: 18
Acq On : 28 Jul 2009 12:54 pm Operator: HC
Sample : 10000ng/ml TO11A Std S21-07270901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:49 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

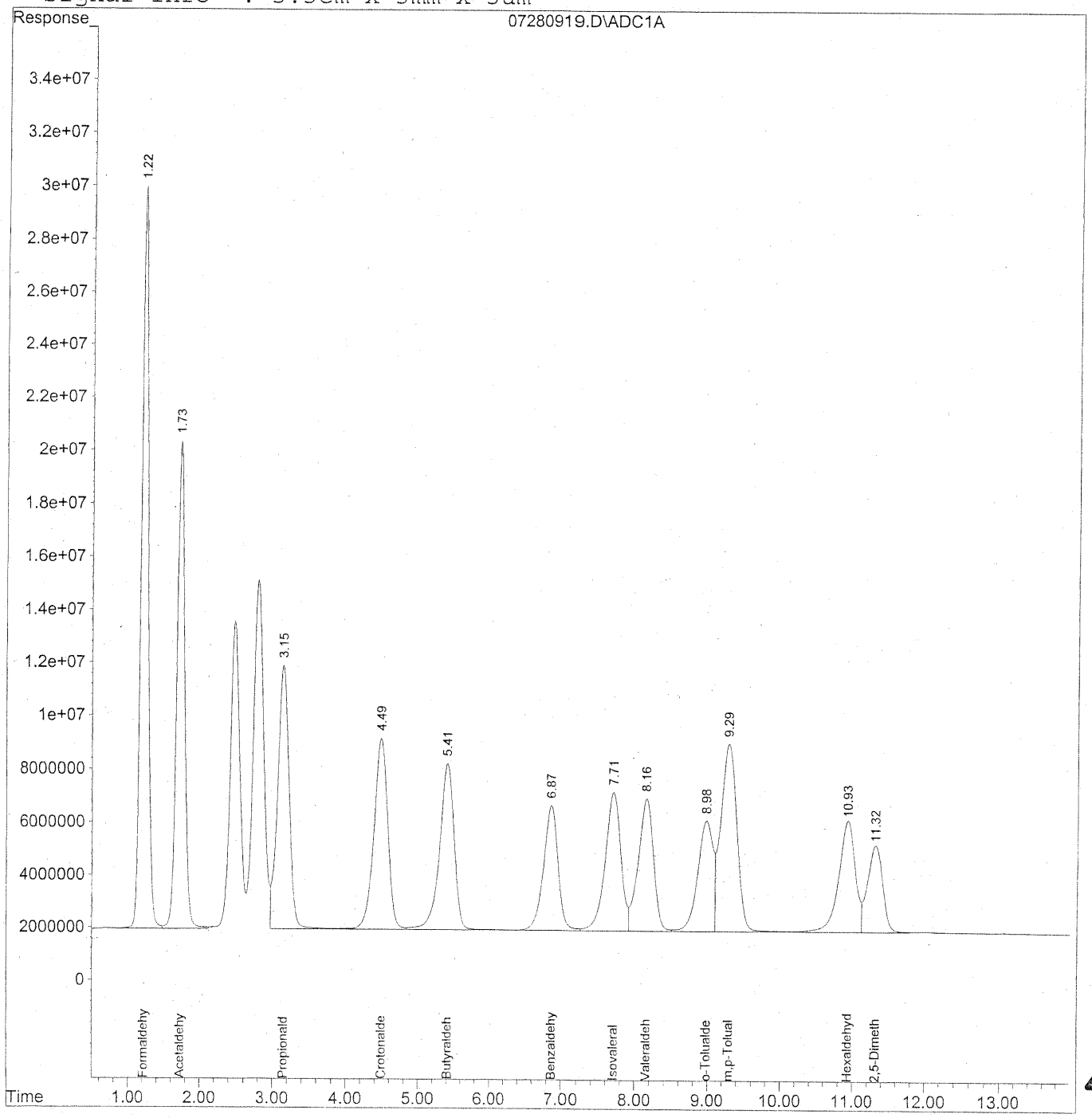
	Compound	R.T.	Response	Conc Units
	Target Compounds			
1)	Formaldehyde	1.22	1908653125	10640.787 ng/ml
2)	Acetaldehyde	1.73	1450154617	10568.343 ng/ml
3)	Propionaldehyde	3.16	1099941045	10538.770 ng/ml
4)	Crotonaldehyde	4.50	972691462	9475.947 ng/ml
5)	Butyraldehyde	5.42	910896701	10789.948 ng/ml
6)	Benzaldehyde	6.88	668462127	10401.737 ng/ml
7)	Isovaleraldehyde	7.72	790328317	9495.299 ng/ml
8)	Valeraldehyde	8.17	730218673	9452.670 ng/ml
9)	o-Tolualdehyde	8.99	608208276	10942.576 ng/ml
10)	m,p-Tolualdehyde	9.30	1111180147	20661.085 ng/ml
11)	Hexaldehyde	10.94	673516807	10163.841 ng/ml
12)	2,5-Dimethylbenzaldehyde	11.32	478460947	9779.540 ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.
07280918.D TO110709.M Tue Jul 28 17:21:16 2009 Page 1

Data File : J:\LC01\DATA\TO11\2009_07\28\07280919.D Vial: 19
Acq On : 28 Jul 2009 1:09 pm Operator: HC
Sample : 10000ng/ml TO11A Std S21-07270901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:50 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



469

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280919.D Vial: 19
Acq On : 28 Jul 2009 1:09 pm Operator: HC
Sample : 10000ng/ml TO11A Std S21-07270901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:50 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	1905913073	10625.511	ng/ml
2)	Acetaldehyde	1.73	1446499891	10541.708	ng/ml
3)	Propionaldehyde	3.15	1098837646	10528.198	ng/ml
4)	Crotonaldehyde	4.49	971357788	9462.954	ng/ml
5)	Butyraldehyde	5.41	911328243	10795.060	ng/ml
6)	Benzaldehyde	6.87	669128969	10412.114	ng/ml
7)	Isovaleraldehyde	7.71	788026190	9467.640	ng/ml
8)	Valeraldehyde	8.16	729839210	9447.758	ng/ml
9)	o-Tolualdehyde	8.98	610326238	10980.681	ng/ml
10)	m,p-Tolualdehyde	9.29	1113209810	20698.824	ng/ml
11)	Hexaldehyde	10.93	681915785	10290.587	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.32	484763918	9908.370	ng/ml

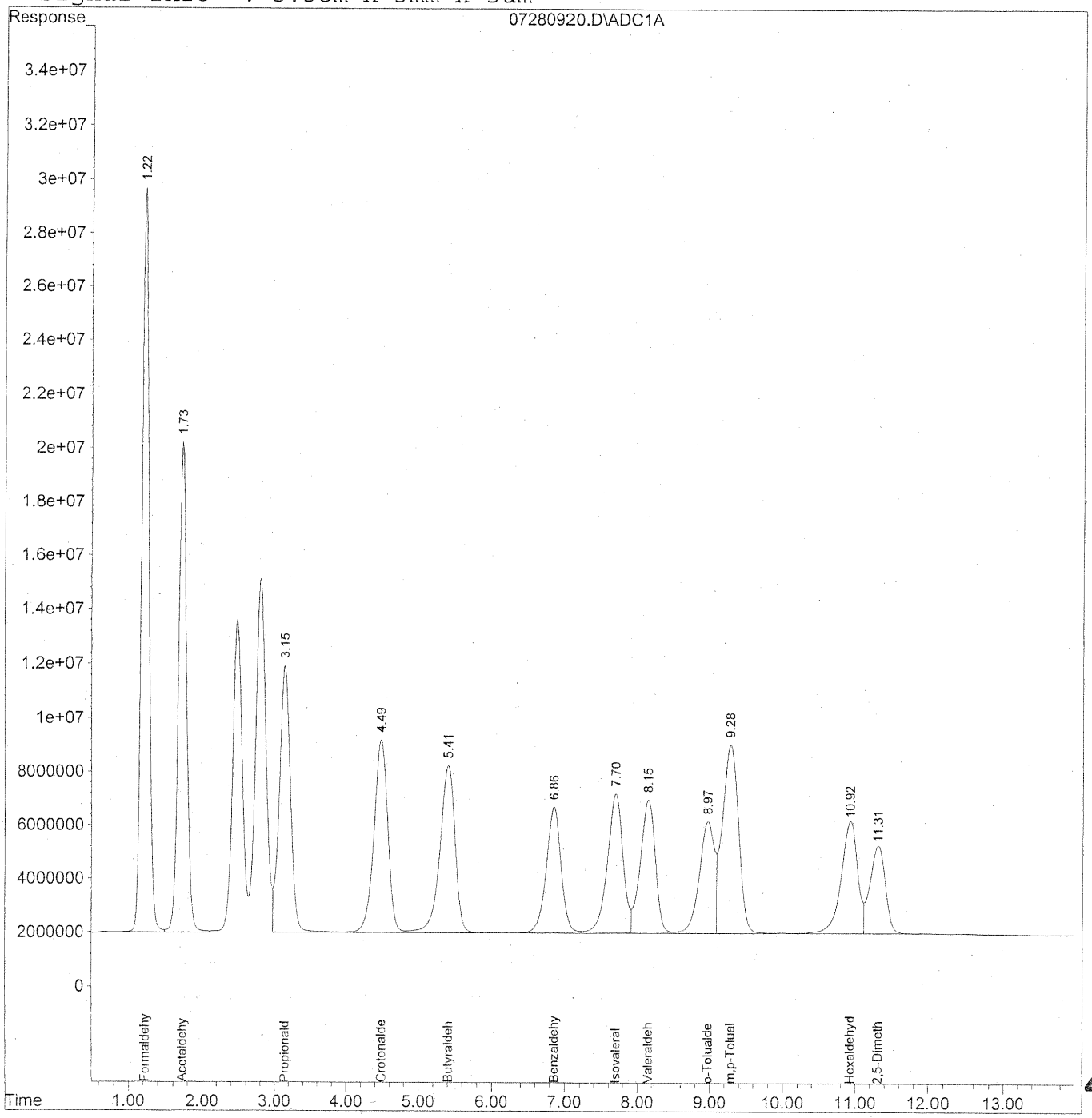
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280920.D Vial: 20
Acq On : 28 Jul 2009 1:25 pm Operator: HC
Sample : 10000ng/ml TO11A Std S21-07270901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:50 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



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Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280920.D Vial: 20
Acq On : 28 Jul 2009 1:25 pm Operator: HC
Sample : 10000ng/ml TO11A Std S21-07270901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:50 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	1875917434	10458.285	ng/ml
2)	Acetaldehyde	1.73	1425028469	10385.230	ng/ml
3)	Propionaldehyde	3.15	1089338811	10437.188	ng/ml
4)	Crotonaldehyde	4.48	963283335	9384.293	ng/ml
5)	Butyraldehyde	5.41	900561239	10667.520	ng/ml
6)	Benzaldehyde	6.86	662238443	10304.892	ng/ml
7)	Isovaleraldehyde	7.70	782256804	9398.325	ng/ml
8)	Valeraldehyde	8.15	722749626	9355.983	ng/ml
9)	o-Tolualdehyde	8.97	603256599	10853.487	ng/ml
10)	m,p-Tolualdehyde	9.29	1100384573	20460.354	ng/ml
11)	Hexaldehyde	10.92	670193360	10113.688	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.31	476113656	9731.563	ng/ml

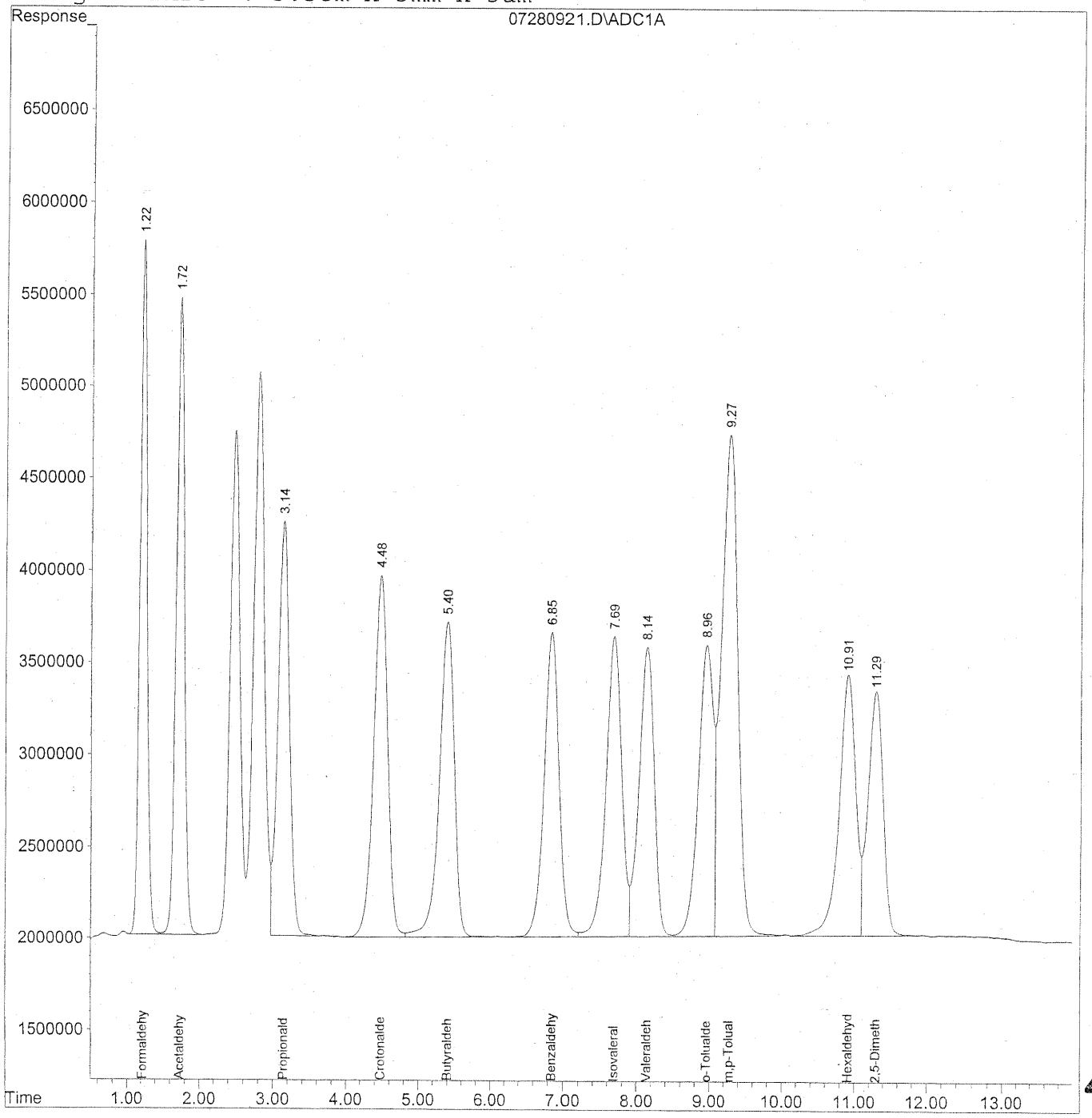
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280921.D Vial: 21
Acq On : 28 Jul 2009 1:40 pm Operator: HC
Sample : ~1500ng/ml TO11A Std ICV S21-07270907 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 17:22 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 15:29:52 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



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Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280921.D Vial: 21
Acq On : 28 Jul 2009 1:40 pm Operator: HC
Sample : ~1500ng/ml TO11A Std ICV S21-07270907 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 17:22 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 15:29:52 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	257076667	1400.342	ng/ml
2)	Acetaldehyde	1.72	270257005	1927.330	ng/ml
3)	Propionaldehyde	3.14	246366252	2309.065	ng/ml
4)	Crotonaldehyde	4.48	262943470	2699.204	ng/ml
5)	Butyraldehyde	5.40	247400524	2800.672	ng/ml
6)	Benzaldehyde	6.85	233067402	3538.331	ng/ml
7)	Isovaleraldehyde	7.69	244473332	3002.720	ng/ml
8)	Valeraldehyde	8.14	226800810	3085.515	ng/ml
9)	o-Tolualdehyde	8.96	225349526	3863.990	ng/ml
10)	m,p-Tolualdehyde	9.27	428359795	7933.265	ng/ml
11)	Hexaldehyde	10.91	226495334	3363.271	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.29	193343187	3944.701	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

TO-11A Aldehyde-DNPH Stock Solution Standard S21-06300801

Source: AccuStandard Inc.
Catalog No: M-8315-R2-DNPH
Lot : B8060121
Solvent: ACN
Expiration Date: 6/12/11

HC 7/29/09

	MW	Aldehyde-DNPH MW*	Manufacturer Prepared Concentration as Aldehyde-DNPH (ug/mL)	Calculated Concentration as Aldehyde (ug/mL)	ICV S21-07270907 (nominal ng/mL)	ICV S21-07270907 (Actual, ng/mL)	% Diff
Formaldehyde	30.03	210.03	100	14.30	1430	1400.34	2.07%
Acetaldehyde	44.05	224.05	100.2	19.70	1970	1927.33	2.17%
Acetone	58.08	238.08	100.2	24.44	2444	not reported	
Acrolein	56.06	236.06	103.1	24.48	2448	not reported	
Propionaldehyde	58.08	238.08	100.2	24.44	2444	2309.07	5.52%
Crotonaldehyde	70.09	250.09	100.2	28.08	2808	2699.20	3.87%
Butyraldehyde	72.11	252.11	100	28.60	2860	2800.67	2.07%
Benzaldehyde	106.12	286.12	100	37.09	3709	3538.33	4.60%
Isovaleraldehyde	86.13	266.13	100.2	32.43	3243	3002.72	7.41%
Valeraldehyde	86.13	266.13	100.1	32.40	3240	3085.52	4.77%
o-Tolualdehyde	120.15	300.15	100.1	40.07	4007	3863.99	3.57%
m,p-Tolualdehyde	120.15	300.15	100.3	80.30	8030	7933.27	1.20%
Hexaldehyde	100.16	280.16	100.3	35.86	3586	3363.27	6.21%
2,5-Dimethylbenzaldehyde	134.18	314.18	100.3	42.84	4284	3944.70	7.92%

(* MW of DNPH is 198g/mol. The result of a nucleophilic reaction of aldehyde & DNPH is a hydrazone derivative with the loss of H20, 18g/mol)

CONTINUING CALIBRATION STANDARDS

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

HC 8/25/09

Printed : 8/25/09
Instrument : LC#1 — Date Acquired : 8/18/09
Detector : UV-VIS 360 — Sample Amount : 5ul
Analyst : HC — Client & PAI Job# : EH&E P0902772

SAMPLE RESULT SUMMARY

Sample Information	MDL	CCV 1500ng/ml S21-08170901	% Diff	ACN CY023 blk	MB front lot 6009/6097 1.0ml	MB back lot 6009/6097 1.0ml	CCV 1500ng/ml S21-08170901	% Diff	P0902772-001 back 1.0ml
Dilution	1.0			1.0	1.0	1.0	1.0		1.0
Sample Volume (L)	NA			NA	NA	NA			0.00
Final Vol.(ml)	1.0			1.0	1.0	1.0	1.0		1.0
	ng/sample	ng/sample		ng/sample	ng/sample	ng/sample	ng/sample		ng/sample
Formaldehyde	100.00	1408.9	6.1%	ND	ND	ND	1407.839	6.1%	ND
Acetaldehyde	100.00	1395.8	6.9%	ND	ND	ND	1394.104	7.1%	ND
Propionaldehyde	100.00	1384.3	7.7%	ND	ND	ND	1388.552	7.4%	ND
Crotonaldehyde	100.00	1364.3	9.0%	ND	ND	ND	1359.517	9.4%	ND
Butyraldehyde	100.00	1415.1	5.7%	ND	ND	ND	1407.627	6.2%	ND
Benzaldehyde	100.00	1364.2	9.1%	ND	ND	ND	1395.367	7.0%	ND
Isovaleraldehyde	100.00	1412.9	5.8%	ND	ND	ND	1422.259	5.2%	ND
Valeraldehyde	100.00	1306.1	12.9%	ND	ND	ND	1334.833	11.0%	ND
o-Tolualdehyde	100.00	1415.6	5.6%	ND	ND	ND	1432.124	4.5%	ND
m,p-Tolualdehyde	200.00	2769.3	7.7%	ND	ND	ND	2762.860	7.9%	ND
Hexaldehyde	100.00	1393.4	7.1%	ND	ND	ND	1421.815	5.2%	ND
2,5-Dimethylbenzaldehyde	100.00	1341.5	10.6%	ND	ND	ND	1309.928	12.7%	ND
	ug/m3			ug/m3	ug/m3	ug/m3			ug/m3
Formaldehyde				NA	NA	NA			ND
Acetaldehyde				NA	NA	NA			ND
Propionaldehyde				NA	NA	NA			ND
Crotonaldehyde				NA	NA	NA			ND
Butyraldehyde				NA	NA	NA			ND
Benzaldehyde				NA	NA	NA			ND
Isovaleraldehyde				NA	NA	NA			ND
Valeraldehyde				NA	NA	NA			ND
o-Tolualdehyde				NA	NA	NA			ND
m,p-Tolualdehyde				NA	NA	NA			ND
Hexaldehyde				NA	NA	NA			ND
2,5-Dimethylbenzaldehyde				NA	NA	NA			ND
	ppb			ppb	ppb	ppb			ppb
Formaldehyde				NA	NA	NA			ND
Acetaldehyde				NA	NA	NA			ND
Propionaldehyde				NA	NA	NA			ND
Crotonaldehyde				NA	NA	NA			ND
Butyraldehyde				NA	NA	NA			ND
Benzaldehyde				NA	NA	NA			ND
Isovaleraldehyde				NA	NA	NA			ND
Valeraldehyde				NA	NA	NA			ND
o-Tolualdehyde				NA	NA	NA			ND
m,p-Tolualdehyde				NA	NA	NA			ND
Hexaldehyde				NA	NA	NA			ND
2,5-Dimethylbenzaldehyde				NA	NA	NA			ND

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

Printed : 8/25/09

Instrument : LC#1 | Date Acquired : 8/18/09

Detector : UV-VIS 360 | Sample Amount : 5ul

Analyst : HC | Client & PAI Job# : EH&E P0902772

HC 8/25/09

SAMPLE RESULT SUMMARY

Sample Information	MDL	P0902772-002 back 1.0ml	P0902772-003 back 1.0ml	P0902772-004 back 1.0ml	P0902772-005 back 1.0ml	P0902772-006 back 1.0ml	CCV 1500ng/ml S21-08170901	% Diff
Dilution	1.0	1.0	1.0	1.0	1.0	1.0	1.0	
Sample Volume (L)	NA	105.60	106.00	109.60	97.60	107.10		
Final Vol.(ml)	1.0	1.0	1.0	1.0	1.0	1.0	1.0	

	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample	
Formaldehyde	100.00	ND	ND	ND	ND	ND	1404.466	6.4%
Acetaldehyde	100.00	403.538 BT	343.535 BT	425.616 BT	ND	283.924 BT	1394.168	7.1%
Propionaldehyde	100.00	ND	~~1882.629 BT~~	ND	ND	ND	1358.521	9.4%
Crotonaldehyde	100.00	ND	ND	ND	ND	ND	1364.214	9.1%
Butyraldehyde	100.00	ND	ND	ND	ND	ND	1405.057	6.3%
Benzaldehyde	100.00	ND	ND	ND	ND	ND	1374.637	8.4%
Isovaleraldehyde	100.00	ND	ND	ND	ND	ND	1428.347	4.8%
Valeraldehyde	100.00	ND	ND	ND	ND	ND	1265.584	15.6%
o-Tolualdehyde	100.00	ND	ND	ND	ND	ND	1429.840	4.7%
m,p-Tolualdehyde	200.00	ND	ND	ND	ND	ND	2769.284	7.7%
Hexaldehyde	100.00	ND	ND	ND	ND	ND	1397.408	6.8%
2,5-Dimethylbenzaldehyde	100.00	ND	ND	ND	ND	ND	1323.025	11.8%

	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3
Formaldehyde		ND	ND	ND	ND	ND
Acetaldehyde		3.821	3.241	3.883	ND	2.651
Propionaldehyde		ND	17.761	ND	ND	ND
Crotonaldehyde		ND	ND	ND	ND	ND
Butyraldehyde		ND	ND	ND	ND	ND
Benzaldehyde		ND	ND	ND	ND	ND
Isovaleraldehyde		ND	ND	ND	ND	ND
Valeraldehyde		ND	ND	ND	ND	ND
o-Tolualdehyde		ND	ND	ND	ND	ND
m,p-Tolualdehyde		ND	ND	ND	ND	ND
Hexaldehyde		ND	ND	ND	ND	ND
2,5-Dimethylbenzaldehyde		ND	ND	ND	ND	ND

	ppb	ppb	ppb	ppb	ppb	ppb
Formaldehyde		ND	ND	ND	ND	ND
Acetaldehyde		2.122	1.800	2.156	ND	1.472
Propionaldehyde		ND	7.480	ND	ND	ND
Crotonaldehyde		ND	ND	ND	ND	ND
Butyraldehyde		ND	ND	ND	ND	ND
Benzaldehyde		ND	ND	ND	ND	ND
Isovaleraldehyde		ND	ND	ND	ND	ND
Valeraldehyde		ND	ND	ND	ND	ND
o-Tolualdehyde		ND	ND	ND	ND	ND
m,p-Tolualdehyde		ND	ND	ND	ND	ND
Hexaldehyde		ND	ND	ND	ND	ND
2,5-Dimethylbenzaldehyde		ND	ND	ND	ND	ND

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

Printed : 8/25/09

Instrument : LC#1 | Date Acquired : 8/18/09

Detector : UV-VIS 360 | Sample Amount : 5ul

Analyst : HC | Client & PAI Job# : EH&E P0902772

Sample Information	MDL	ACN blk lot CY023	MB front lot 6009/6097 1.0ml	MB back lot 6009/6097 1.0ml	P0902772-007 back 1.0ml	P0902772-008 back 1.0ml	P0902772-009 back1.0ml	P0902772-010 back 1.0ml
Dilution	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
Sample Volume (L)	NA				106.30	108.20	0.00	110.90
Final Vol.(ml)	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample
Formaldehyde	100.00	ND	ND	ND	ND	ND	ND	ND
Acetaldehyde	100.00	ND	ND	ND	1865.973	1770.389	ND	ND
Propionaldehyde	100.00	ND	ND	ND	ND	ND	ND	ND
Crotonaldehyde	100.00	ND	ND	ND	ND	ND	ND	ND
Butyraldehyde	100.00	ND	ND	ND	ND	ND	ND	ND
Benzaldehyde	100.00	ND	ND	ND	ND	ND	ND	ND
Isovaleraldehyde	100.00	ND	ND	ND	ND	ND	ND	ND
Valeraldehyde	100.00	ND	ND	ND	ND	ND	ND	ND
o-Tolualdehyde	100.00	ND	ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde	200.00	ND	ND	ND	ND	ND	ND	ND
Hexaldehyde	100.00	ND	ND	ND	ND	ND	ND	ND
2,5-Dimethylbenzaldehyde	100.00	ND	ND	ND	ND	ND	ND	ND
	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3
Formaldehyde		ND	ND	ND	ND	ND	ND	ND
Acetaldehyde		ND	ND	ND	17.554	16.362	ND	ND
Propionaldehyde		ND	ND	ND	ND	ND	ND	ND
Crotonaldehyde		ND	ND	ND	ND	ND	ND	ND
Butyraldehyde		ND	ND	ND	ND	ND	ND	ND
Benzaldehyde		ND	ND	ND	ND	ND	ND	ND
Isovaleraldehyde		ND	ND	ND	ND	ND	ND	ND
Valeraldehyde		ND	ND	ND	ND	ND	ND	ND
o-Tolualdehyde		ND	ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde		ND	ND	ND	ND	ND	ND	ND
Hexaldehyde		ND	ND	ND	ND	ND	ND	ND
2,5-Dimethylbenzaldehyde		ND	ND	ND	ND	ND	ND	ND
	ppb	ppb	ppb	ppb	ppb	ppb	ppb	ppb
Formaldehyde		ND	ND	ND	ND	ND	ND	ND
Acetaldehyde		ND	ND	ND	9.747	9.086	ND	ND
Propionaldehyde		ND	ND	ND	ND	ND	ND	ND
Crotonaldehyde		ND	ND	ND	ND	ND	ND	ND
Butyraldehyde		ND	ND	ND	ND	ND	ND	ND
Benzaldehyde		ND	ND	ND	ND	ND	ND	ND
Isovaleraldehyde		ND	ND	ND	ND	ND	ND	ND
Valeraldehyde		ND	ND	ND	ND	ND	ND	ND
o-Tolualdehyde		ND	ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde		ND	ND	ND	ND	ND	ND	ND
Hexaldehyde		ND	ND	ND	ND	ND	ND	ND
2,5-Dimethylbenzaldehyde		ND	ND	ND	ND	ND	ND	ND

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

		Printed :	8/25/09
Instrument :	LC#1	Date Acquired :	8/18/09
Detector :	UV-VIS 360	Sample Amount 5ul	
Analyst :	HC	Client & PAI Job EH&E P0902772	

ll 8/25/09

Sample Information	MDL	P0902772-011 back 1.0ml	P0902772-012 back 1.0ml	CCV 1500ng/ml S21-08170902	% Diff	ACN Blk lot CY023	MB front lot 6009/6097 1.0ml	MB back lot 6009/6097 1.0ml
Dilution	1.0	1.0	1.0	1.0		1.0	1.0	1.0
Sample Volume (L)	NA	100.60	103.70					
Final Vol.(ml)	1.0	1.0	1.0	1.0		1.0	1.0	1.0

	ng/sample	ng/sample	ng/sample	ng/sample		ng/sample	ng/sample	ng/sample
Formaldehyde	100.00	ND	ND	1408.550	6.1%	ND	ND	ND
Acetaldehyde	100.00	1814.948 BT	1795.262 BT	1391.222	7.3%	ND	ND	ND
Propionaldehyde	100.00	ND	ND	1374.156	8.4%	ND	ND	ND
Crotonaldehyde	100.00	ND	ND	1366.856	8.9%	ND	ND	ND
Butyraldehyde	100.00	ND	ND	1403.759	6.4%	ND	ND	ND
Benzaldehyde	100.00	ND	ND	1391.194	7.3%	ND	ND	ND
Isovaleraldehyde	100.00	ND	ND	1456.549	2.9%	ND	ND	ND
Valeraldehyde	100.00	ND	ND	1300.659	13.3%	ND	ND	ND
o-Tolualdehyde	100.00	ND	ND	1421.315	5.2%	ND	ND	ND
m,p-Tolualdehyde	200.00	ND	ND	2790.978	7.0%	ND	ND	ND
Hexaldehyde	100.00	ND	ND	1392.264	7.2%	ND	ND	ND
2,5-Dimethylbenzaldehyde	100.00	ND	ND	1318.842	12.1%	ND	ND	ND

	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3
Formaldehyde	ND	ND		ND	ND	ND
Acetaldehyde	18.041	17.312		ND	ND	ND
Propionaldehyde	ND	ND		ND	ND	ND
Crotonaldehyde	ND	ND		ND	ND	ND
Butyraldehyde	ND	ND		ND	ND	ND
Benzaldehyde	ND	ND		ND	ND	ND
Isovaleraldehyde	ND	ND		ND	ND	ND
Valeraldehyde	ND	ND		ND	ND	ND
o-Tolualdehyde	ND	ND		ND	ND	ND
m,p-Tolualdehyde	ND	ND		ND	ND	ND
Hexaldehyde	ND	ND		ND	ND	ND
2,5-Dimethylbenzaldehyde	ND	ND		ND	ND	ND

	ppb	ppb	ppb	ppb	ppb	ppb
Formaldehyde	ND	ND		ND	ND	ND
Acetaldehyde	10.018	9.613		ND	ND	ND
Propionaldehyde	ND	ND		ND	ND	ND
Crotonaldehyde	ND	ND		ND	ND	ND
Butyraldehyde	ND	ND		ND	ND	ND
Benzaldehyde	ND	ND		ND	ND	ND
Isovaleraldehyde	ND	ND		ND	ND	ND
Valeraldehyde	ND	ND		ND	ND	ND
o-Tolualdehyde	ND	ND		ND	ND	ND
m,p-Tolualdehyde	ND	ND		ND	ND	ND
Hexaldehyde	ND	ND		ND	ND	ND
2,5-Dimethylbenzaldehyde	ND	ND		ND	ND	ND

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

		Printed :	8/25/09
Instrument :	LC#1	Date Acquirec	8/18/09
Detector :	UV-VIS 360	Sample Amou	5ul
Analyst :	HC	Client & PAI	J EH&E P0902772

tec 8/25/09

Sample Information	MDL	CCV 1500ng/ml S21-08170902	% Diff	P0902772-001 front 1.0ml	P0902772-002 front 1.0ml	P0902772-003 front 1.0ml	P0902772-004 front 1.0ml	P0902772-005 front 1.0ml	P0902772-006 front 1.0ml
Dilution	1.0	1.0		1.0	1.0	1.0	1.0	1.0	1.0
Sample Volume (L)	NA			0.00	105.60	106.00	109.60	97.60	107.10
Final Vol.(ml)	1.0	1.0		1.0	1.0	1.0	1.0	1.0	1.0
	ng/sample	ng/sample		ng/sample	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample
Formaldehyde	100.00	1386.265	7.6%	ND	3842.366	3720.357	4156.856	ND	3991.607
Acetaldehyde	100.00	1374.461	8.4%	ND	2181.480	2223.205	2237.807	ND	2305.320
Propionaldehyde	100.00	1342.093	10.5%	ND	197.478	219.598	220.236	ND	233.132
Crotonaldehyde	100.00	1342.545	10.5%	ND	ND	ND	ND	ND	ND
Butyraldehyde	100.00	1394.032	7.1%	ND	209.888	223.368	240.626	ND	426.423
Benzaldehyde	100.00	1386.419	7.6%	ND	313.417	344.570	352.343	ND	409.326
Isovaleraldehyde	100.00	1382.381	7.8%	ND	ND	ND	ND	ND	ND
Valeraldehyde	100.00	1307.223	12.9%	ND	245.917	238.403	332.014	ND	380.917
o-Tolualdehyde	100.00	1395.485	7.0%	ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde	200.00	2726.214	9.1%	ND	ND	ND	ND	ND	ND
Hexaldehyde	100.00	1369.516	8.7%	ND	612.936	781.477	669.492	ND	624.050
2,5-Dimethylbenzaldehyde	100.00	1283.584	14.4%	ND	489.222	744.633	577.570	731.970	547.292

ug/m3		ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3
Formaldehyde		ND	36.386	35.098	37.928	ND	37.270
Acetaldehyde		ND	20.658	20.974	20.418	ND	21.525
Propionaldehyde		ND	1.870	2.072	2.009	ND	2.177
Crotonaldehyde		ND	ND	ND	ND	ND	ND
Butyraldehyde		ND	1.988	2.107	2.195	ND	3.982
Benzaldehyde		ND	2.968	3.251	3.215	ND	3.822
Isovaleraldehyde		ND	ND	ND	ND	ND	ND
Valeraldehyde		ND	2.329	2.249	3.029	ND	3.557
o-Tolualdehyde		ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde		ND	ND	ND	ND	ND	ND
Hexaldehyde		ND	5.804	7.372	6.109	ND	5.827
2,5-Dimethylbenzaldehyde		ND	4.633	7.025	5.270	7.500	5.110

ppb		ppb	ppb	ppb	ppb	ppb	ppb
Formaldehyde		ND	29.637	28.588	30.893	ND	30.357
Acetaldehyde		ND	11.471	11.646	11.338	ND	11.952
Propionaldehyde		ND	0.788	0.872	0.846	ND	0.917
Crotonaldehyde		ND	ND	ND	ND	ND	ND
Butyraldehyde		ND	0.674	0.715	0.745	ND	1.351
Benzaldehyde		ND	0.684	0.749	0.741	ND	0.881
Isovaleraldehyde		ND	ND	ND	ND	ND	ND
Valeraldehyde		ND	0.661	0.639	0.860	ND	1.010
o-Tolualdehyde		ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde		ND	ND	ND	ND	ND	ND
Hexaldehyde		ND	1.417	1.800	1.492	ND	1.423
2,5-Dimethylbenzaldehyde		ND	0.845	1.281	0.961	1.367	0.932

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

Printed : 8/25/09
Instrument : LC#1
Date Acquirec 8/18/09
Detector : UV-VIS 360
Sample Amou 5ul
Analyst : HC
Client & PAI J EH&E P0902772

HC 8/25/09

Sample Information	MDL	CCV 1500ng/ml S21-08170902	% Diff	ACN blk lot CY023	MB front lot 6009/6097 1.0ml	MB back lot 6009/6097 1.0ml	P0902772-007 front 1.0ml	P0902772-008 front 1.0ml	P0902772-009 front 1.0ml
Dilution	1.0	1.0		1.0	1.0	1.0	1.0	1.0	1.0
Sample Volume (L)	NA						106.30	108.20	0.00
Final Vol.(ml)	1.0	1.0		1.0	1.0	1.0	1.0	1.0	1.0

	ng/sample	ng/sample		ng/sample	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample
Formaldehyde	100.00	1448.873	3.4%	ND	ND	ND	~~12705.410~~	~~12405.985~~ SD	ND
Acetaldehyde	100.00	1445.682	3.6%	ND	ND	ND	~~13138.620~~	~~13678.089~~ SD	ND
Propionaldehyde	100.00	1412.476	5.8%	ND	ND	ND	876.671	881.332	ND
Crotonaldehyde	100.00	1387.219	7.5%	ND	ND	ND	ND	ND	ND
Butyraldehyde	100.00	1429.351	4.7%	ND	ND	ND	527.459	537.335	ND
Benzaldehyde	100.00	1413.272	5.8%	ND	ND	ND	1038.895	1182.769	ND
Isovaleraldehyde	100.00	1454.238	3.1%	ND	ND	ND	277.520	393.376	ND
Valeraldehyde	100.00	1346.625	10.2%	ND	ND	ND	1050.376	1448.289	ND
o-Tolualdehyde	100.00	1464.162	2.4%	ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde	200.00	2856.189	4.8%	ND	ND	ND	ND	ND	ND
Hexaldehyde	100.00	1385.281	7.6%	ND	ND	ND	2655.258	3082.279	ND
2,5-Dimethylbenzaldehyde	100.00	1313.238	12.5%	ND	ND	ND	ND	230.567	ND

ug/m3		ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3
Formaldehyde		ND	ND	ND	119.524	114.658	ND
Acetaldehyde		ND	ND	ND	123.599	126.415	ND
Propionaldehyde		ND	ND	ND	8.247	8.145	ND
Crotonaldehyde		ND	ND	ND	ND	ND	ND
Butyraldehyde		ND	ND	ND	4.962	4.966	ND
Benzaldehyde		ND	ND	ND	9.773	10.931	ND
Isovaleraldehyde		ND	ND	ND	2.611	3.636	ND
Valeraldehyde		ND	ND	ND	9.881	13.385	ND
o-Tolualdehyde		ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde		ND	ND	ND	ND	ND	ND
Hexaldehyde		ND	ND	ND	24.979	28.487	ND
2,5-Dimethylbenzaldehyde		ND	ND	ND	ND	2.131	ND

ppb		ppb	ppb	ppb	ppb	ppb	ppb
Formaldehyde		ND	ND	ND	97.355	93.391	ND
Acetaldehyde		ND	ND	ND	68.632	70.195	ND
Propionaldehyde		ND	ND	ND	3.473	3.430	ND
Crotonaldehyde		ND	ND	ND	ND	ND	ND
Butyraldehyde		ND	ND	ND	1.683	1.685	ND
Benzaldehyde		ND	ND	ND	2.253	2.520	ND
Isovaleraldehyde		ND	ND	ND	0.741	1.032	ND
Valeraldehyde		ND	ND	ND	2.806	3.801	ND
o-Tolualdehyde		ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde		ND	ND	ND	ND	ND	ND
Hexaldehyde		ND	ND	ND	6.100	6.957	ND
2,5-Dimethylbenzaldehyde		ND	ND	ND	ND	0.388	ND

SD = See dilution HC 8/25/09

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

Printed : 8/25/09
Instrument : LC#1
Date Acquirec 8/18/09
Detector : UV-VIS 360
Sample Amou 5ul
Analyst : HC
Client & PAI J EH&E P0902772

HC 8/25/09

Sample Information	MDL	P0902772-010 front 1.0ml	P0902772-011 front 1.0ml	P0902772-012 front 1.0ml	CCV 1500ng/ml S21-08180901	% Diff	ACN lot CY023	P0902772-007 front 10x	P0902772-008 front 10x
Dilution	1.0	1.0	1.0	1.0	1.0		1.0	10.0	10.0
Sample Volume (L)	NA	110.90	100.60	103.70				106.30	108.20
Final Vol.(ml)	1.0	1.0	1.0	1.0	1.0		1.0	1.0	1.0
	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample		ng/sample	ng/sample	ng/sample
Formaldehyde	100.00	410.006	~~11574.292~~ SD	~~12014.688~~ SD	1421.628	5.2%	ND	12107.530	12107.420
Acetaldehyde	100.00	191.241	~~12084.362~~ SD	~~12319.756~~ SD	1392.656	7.2%	ND	12817.000	13553.490
Propionaldehyde	100.00	ND	788.850	863.185	1392.089	7.2%	ND		
Crotonaldehyde	100.00	ND	ND	ND	1361.969	9.2%	ND		
Butyraldehyde	100.00	ND	478.162	447.746	1391.429	7.2%	ND		
Benzaldehyde	100.00	ND	1121.396	1337.327	1366.266	8.9%	ND		
Isovaleraldehyde	100.00	ND	237.056	339.469	1399.726	6.7%	ND		
Valeraldehyde	100.00	ND	897.205	1163.409	1322.256	11.8%	ND		
o-Tolualdehyde	100.00	ND	ND	ND	1410.584	6.0%	ND		
m,p-Tolualdehyde	200.00	ND	ND	ND	2797.883	6.7%	ND		
Hexaldehyde	100.00	ND	2658.393	~~2989.804~~	1428.555	4.8%	ND		
2,5-Dimethylbenzaldehyde	100.00	ND	ND	ND	1310.223	12.7%	ND		
	ug/m3	ug/m3	ug/m3	ug/m3			ug/m3	ug/m3	ug/m3
Formaldehyde		3.697	115.053	115.860			ND	113.900	111.899
Acetaldehyde		1.724	120.123	118.802			ND	120.574	125.263
Propionaldehyde		ND	7.841	8.324			ND		
Crotonaldehyde		ND	ND	ND			ND		
Butyraldehyde		ND	4.753	4.318			ND		
Benzaldehyde		ND	11.147	12.896			ND		
Isovaleraldehyde		ND	2.356	3.274			ND		
Valeraldehyde		ND	8.919	11.219			ND		
o-Tolualdehyde		ND	ND	ND			ND		
m,p-Tolualdehyde		ND	ND	ND			ND		
Hexaldehyde		ND	26.425	28.831			ND		
2,5-Dimethylbenzaldehyde		ND	ND	ND			ND		
	ppb	ppb	ppb	ppb			ppb	ppb	ppb
Formaldehyde		3.011	93.713	94.370			ND	92.773	91.143
Acetaldehyde		0.958	66.702	65.968			ND	66.952	69.556
Propionaldehyde		ND	3.302	3.506			ND		
Crotonaldehyde		ND	ND	ND			ND		
Butyraldehyde		ND	1.612	1.465			ND		
Benzaldehyde		ND	2.569	2.972			ND		
Isovaleraldehyde		ND	0.669	0.930			ND		
Valeraldehyde		ND	2.533	3.186			ND		
o-Tolualdehyde		ND	ND	ND			ND		
m,p-Tolualdehyde		ND	ND	ND			ND		
Hexaldehyde		ND	6.453	7.041			ND		
2,5-Dimethylbenzaldehyde		ND	ND	ND			ND		

SD = see dilution HC 8/25/09

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

		Printed :	8/25/09
Instrument :	LC#1	Date Acquirec	8/18/09
Detector :	UV-VIS 360	Sample Amou	5ul
Analyst :	HC	Client & PAI	J EH&E P0902772

HC 8/25/09

Sample Information	MDL	P0902772-011 front 10x	P0902772-012 front 10x	CCV 1500ng/ml S21-08180901	% Diff
Dilution	1.0	10.0	10.0	1.0	
Sample Volume (L)	NA	100.60	103.70		
Final Vol.(ml)	1.0	1.0	1.0	1.0	
	ng/sample	ng/sample	ng/sample	ng/sample	
Formaldehyde	100.00	11229.930	11536.130	1413.127	5.8%
Acetaldehyde	100.00	12036.670	12106.650	1389.691	7.4%
Propionaldehyde	100.00			1387.048	7.5%
Crotonaldehyde	100.00			1355.494	9.6%
Butyraldehyde	100.00			1416.651	5.6%
Benzaldehyde	100.00			1393.704	7.1%
Isovaleraldehyde	100.00			1408.378	6.1%
Valeraldehyde	100.00			1367.744	8.8%
o-Tolualdehyde	100.00			1433.404	4.4%
m,p-Tolualdehyde	200.00			2812.368	6.3%
Hexaldehyde	100.00			1507.739	0.5%
2,5-Dimethylbenzaldehyde	100.00			1420.073	5.3%

	ug/m3	ug/m3	ug/m3
Formaldehyde		111.630	111.245
Acetaldehyde		119.649	116.747
Propionaldehyde			
Crotonaldehyde			
Butyraldehyde			
Benzaldehyde			
Isovaleraldehyde			
Valeraldehyde			
o-Tolualdehyde			
m,p-Tolualdehyde			
Hexaldehyde			
2,5-Dimethylbenzaldehyde			

	ppb	ppb	ppb
Formaldehyde		90.924	90.611
Acetaldehyde		66.438	64.827
Propionaldehyde			
Crotonaldehyde			
Butyraldehyde			
Benzaldehyde			
Isovaleraldehyde			
Valeraldehyde			
o-Tolualdehyde			
m,p-Tolualdehyde			
Hexaldehyde			
2,5-Dimethylbenzaldehyde			

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

Printed : 8/25/09
Instrument : LC#1
Date Acquired : 8/17/09
Detector : UV-VIS 360
Sample Amount : 5ul
Analyst : HC
Client & PAI Job# : EH&E P0902772

SAMPLE RESULT SUMMARY

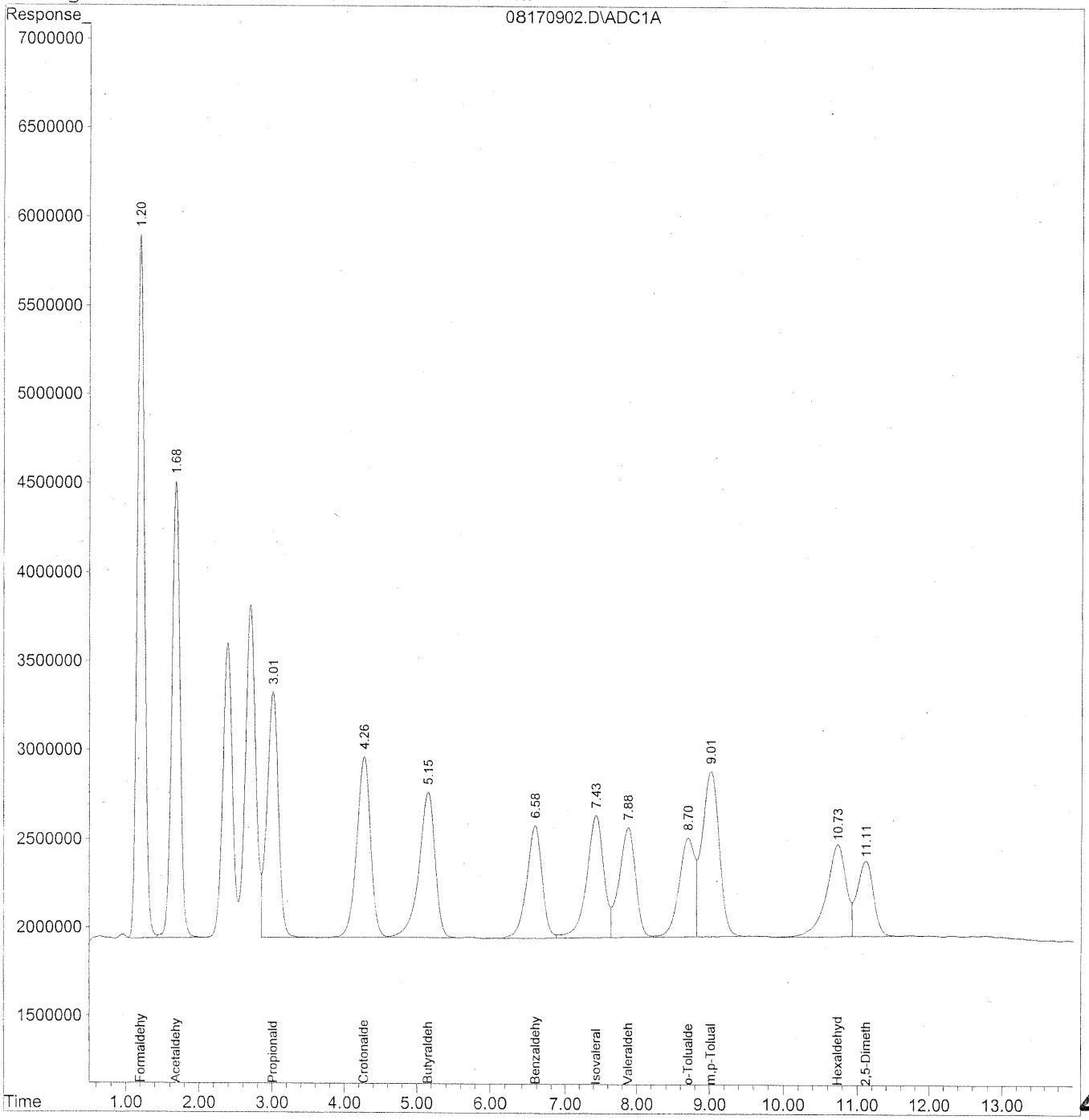
Sample Information	MDL	CCV 1500ng/ml S21-08170901	% Diff	CCV 1500ng/ml S21-08170902	% Diff
Dilution	1.0			1.0	
Sample Volume (L)	NA			NA	
Final Vol.(ml)	1.0			1.0	
	ng/sample	ng/sample		ng/sample	
Formaldehyde	100.00	1415.1	5.7%	1426.460	4.9%
Acetaldehyde	100.00	1388.5	7.4%	1417.099	5.5%
Propionaldehyde	100.00	1382.4	7.8%	1392.524	7.2%
Crotonaldehyde	100.00	1360.1	9.3%	1373.100	8.5%
Butyraldehyde	100.00	1404.5	6.4%	1423.311	5.1%
Benzaldehyde	100.00	1406.1	6.3%	1405.775	6.3%
Isovaleraldehyde	100.00	1476.6	1.6%	1434.947	4.3%
Valeraldehyde	100.00	1295.7	13.6%	1319.212	12.1%
o-Tolualdehyde	100.00	1439.0	4.1%	1434.876	4.3%
m,p-Tolualdehyde	200.00	2794.7	6.8%	2792.447	6.9%
Hexaldehyde	100.00	1425.3	5.0%	1410.245	6.0%
2,5-Dimethylbenzaldehyde	100.00	1317.7	12.2%	1297.199	13.5%
	ug/m3			ug/m3	ug/m3
Formaldehyde				NA	NA
Acetaldehyde				NA	NA
Propionaldehyde				NA	NA
Crotonaldehyde				NA	NA
Butyraldehyde				NA	NA
Benzaldehyde				NA	NA
Isovaleraldehyde				NA	NA
Valeraldehyde				NA	NA
o-Tolualdehyde				NA	NA
m,p-Tolualdehyde				NA	NA
Hexaldehyde				NA	NA
2,5-Dimethylbenzaldehyde				NA	NA
	ppb			ppb	ppb
Formaldehyde				NA	NA
Acetaldehyde				NA	NA
Propionaldehyde				NA	NA
Crotonaldehyde				NA	NA
Butyraldehyde				NA	NA
Benzaldehyde				NA	NA
Isovaleraldehyde				NA	NA
Valeraldehyde				NA	NA
o-Tolualdehyde				NA	NA
m,p-Tolualdehyde				NA	NA
Hexaldehyde				NA	NA
2,5-Dimethylbenzaldehyde				NA	NA

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170902.D Vial: 2
Acq On : 17 Aug 2009 3:05 pm Operator: HC
Sample : 1500ng/ml TO11A std S21-08170901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:04 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170902.D Vial: 2
Acq On : 17 Aug 2009 3:05 pm Operator: HC
Sample : 1500ng/ml TO11A std S21-08170901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:04 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

HC 8/19/09

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.20	266825332	1453.444	ng/ml
2)	Acetaldehyde	1.68	202093866	1441.227	ng/ml
3)	Propionaldehyde	3.01	151971655	1424.353	ng/ml
4)	Crotonaldehyde	4.26	134978482	1385.600	ng/ml
5)	Butyraldehyde	5.15	127063547	1438.410	ng/ml
6)	Benzaldehyde	6.57	92950843	1411.140	ng/ml
7)	Isovaleraldehyde	7.43	115183219	1471.971	ng/ml
8)	Valeraldehyde	7.88	97937916	1332.398	ng/ml
9)	o-Tolualdehyde	8.70	83932300	1439.158	ng/ml
10)	m,p-Tolualdehyde	9.02	153037610	2834.271	ng/ml
11)	Hexaldehyde	10.73	96954329	1439.693	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.11	66414722	1355.032	ng/ml

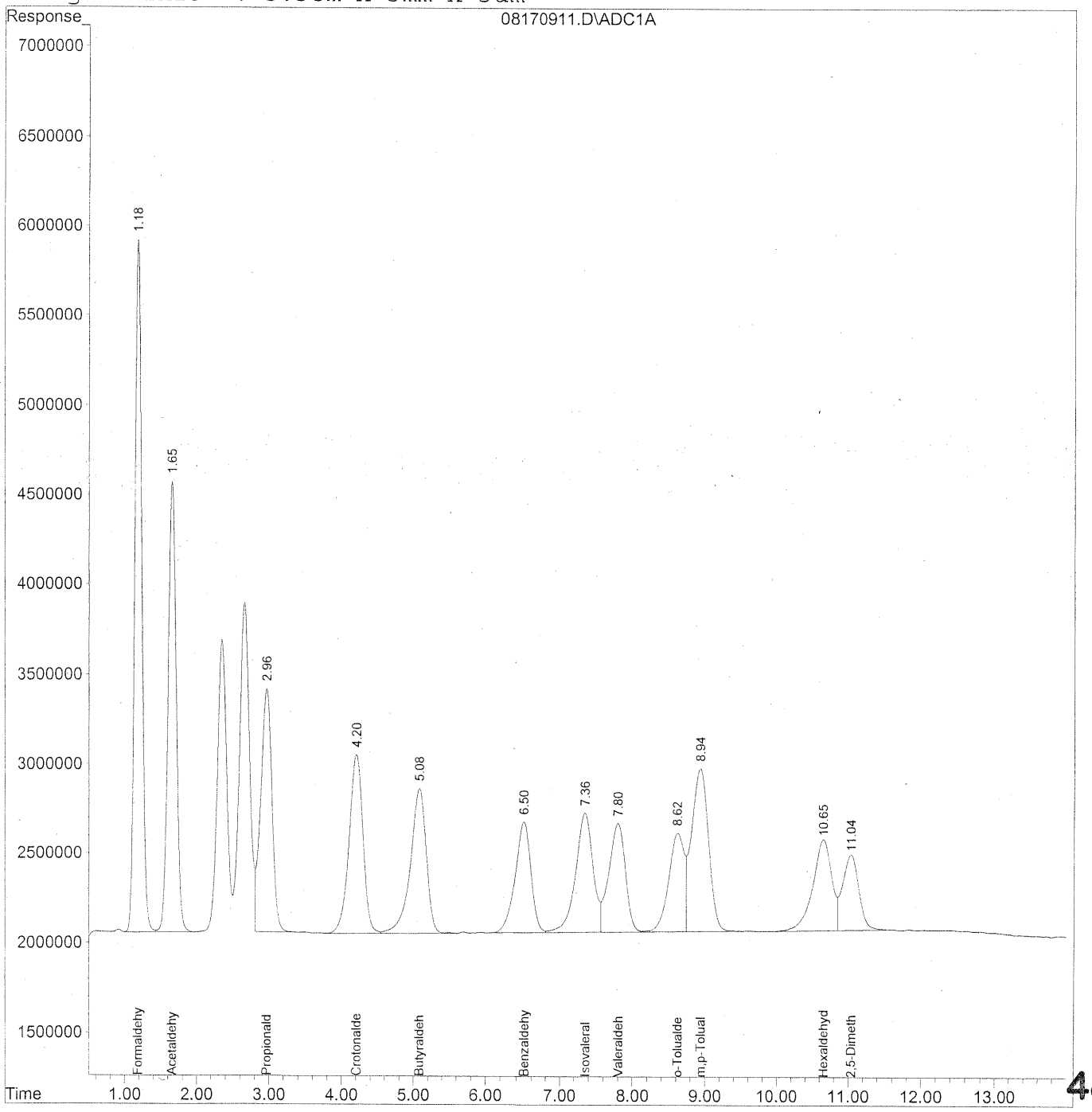
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170911.D Vial: 11
Acq On : 17 Aug 2009 5:20 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 17 18:21 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 17 11:19:09 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



488

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170911.D Vial: 11
Acq On : 17 Aug 2009 5:20 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 17 18:21 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 17 11:19:09 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
Target Compounds					
1)	Formaldehyde	1.18	258649524	1408.909	ng/ml
2)	Acetaldehyde	1.65	195724864	1395.806	ng/ml
3)	Propionaldehyde	2.96	147694855	1384.269	ng/ml
4)	Crotonaldehyde	4.20	132908163	1364.348	ng/ml
5)	Butyraldehyde	5.08	125004364	1415.099	ng/ml
6)	Benzaldehyde	6.50	89857092	1364.172	ng/ml
7)	Isovaleraldehyde	7.35	110564037	1412.941	ng/ml
8)	Valeraldehyde	7.80	96007837	1306.140	ng/ml
9)	o-Tolualdehyde	8.62	82559088	1415.612	ng/ml
10)	m,p-Tolualdehyde	8.94	149526793	2769.251	ng/ml
11)	Hexaldehyde	10.65	93835328	1393.378	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.04	65751013	1341.491	ng/ml

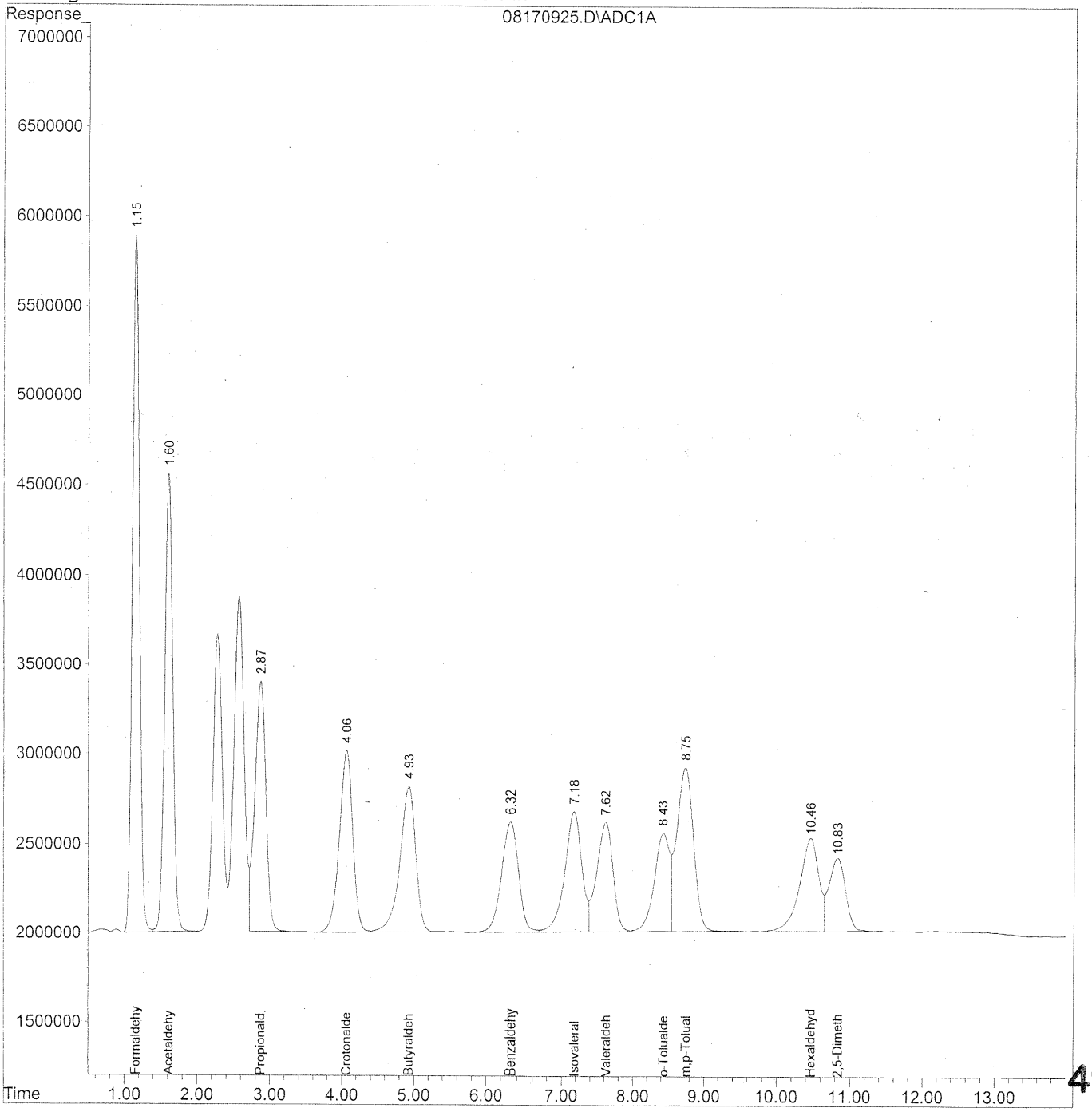
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170925.D Vial: 25
Acq On : 17 Aug 2009 8:51 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 18 8:05 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 17 18:29:01 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



490

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170925.D Vial: 25
Acq On : 17 Aug 2009 8:51 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 18 8:05 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 17 18:29:01 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

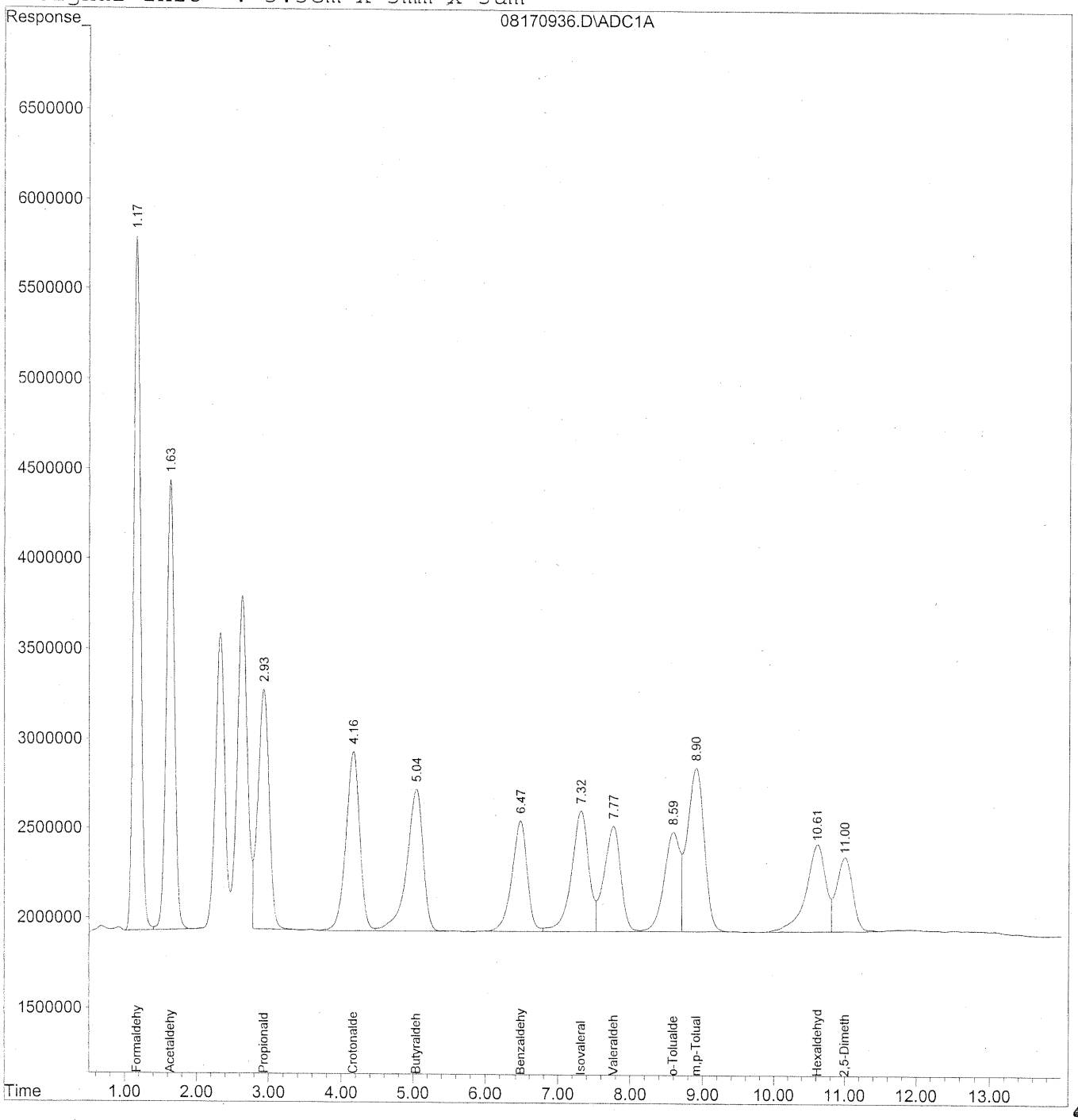
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.15	258452988	1407.839	ng/ml
2)	Acetaldehyde	1.61	195486130	1394.104	ng/ml
3)	Propionaldehyde	2.87	148151844	1388.552	ng/ml
4)	Crotonaldehyde	4.06	132437596	1359.517	ng/ml
5)	Butyraldehyde	4.93	124344289	1407.627	ng/ml
6)	Benzaldehyde	6.32	91911870	1395.367	ng/ml
7)	Isovaleraldehyde	7.18	111293183	1422.259	ng/ml
8)	Valeraldehyde	7.62	98116879	1334.833	ng/ml
9)	o-Tolualdehyde	8.43	83522070	1432.124	ng/ml
10)	m,p-Tolualdehyde	8.74	149181730	2762.860	ng/ml
11)	Hexaldehyde	10.46	95750360	1421.815	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.83	64204013	1309.928	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Data File : J:\LC01\DATA\TO11\2009_08\17\08170936.D Vial: 35
Acq On : 17 Aug 2009 11:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:02 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170936.D Vial: 35
Acq On : 17 Aug 2009 11:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:02 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	257833865	1404.466	ng/ml
2)	Acetaldehyde	1.63	195495143	1394.168	ng/ml
3)	Propionaldehyde	2.93	144947661	1358.521	ng/ml
4)	Crotonaldehyde	4.16	132895137	1364.214	ng/ml
5)	Butyraldehyde	5.04	124117271	1405.057	ng/ml
6)	Benzaldehyde	6.46	90546374	1374.637	ng/ml
7)	Isovaleraldehyde	7.32	111769555	1428.347	ng/ml
8)	Valeraldehyde	7.77	93026780	1265.584	ng/ml
9)	o-Tolualdehyde	8.59	83388888	1429.840	ng/ml
10)	m,p-Tolualdehyde	8.91	149528624	2769.284	ng/ml
11)	Hexaldehyde	10.61	94106711	1397.408	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.00	64845936	1323.025	ng/ml

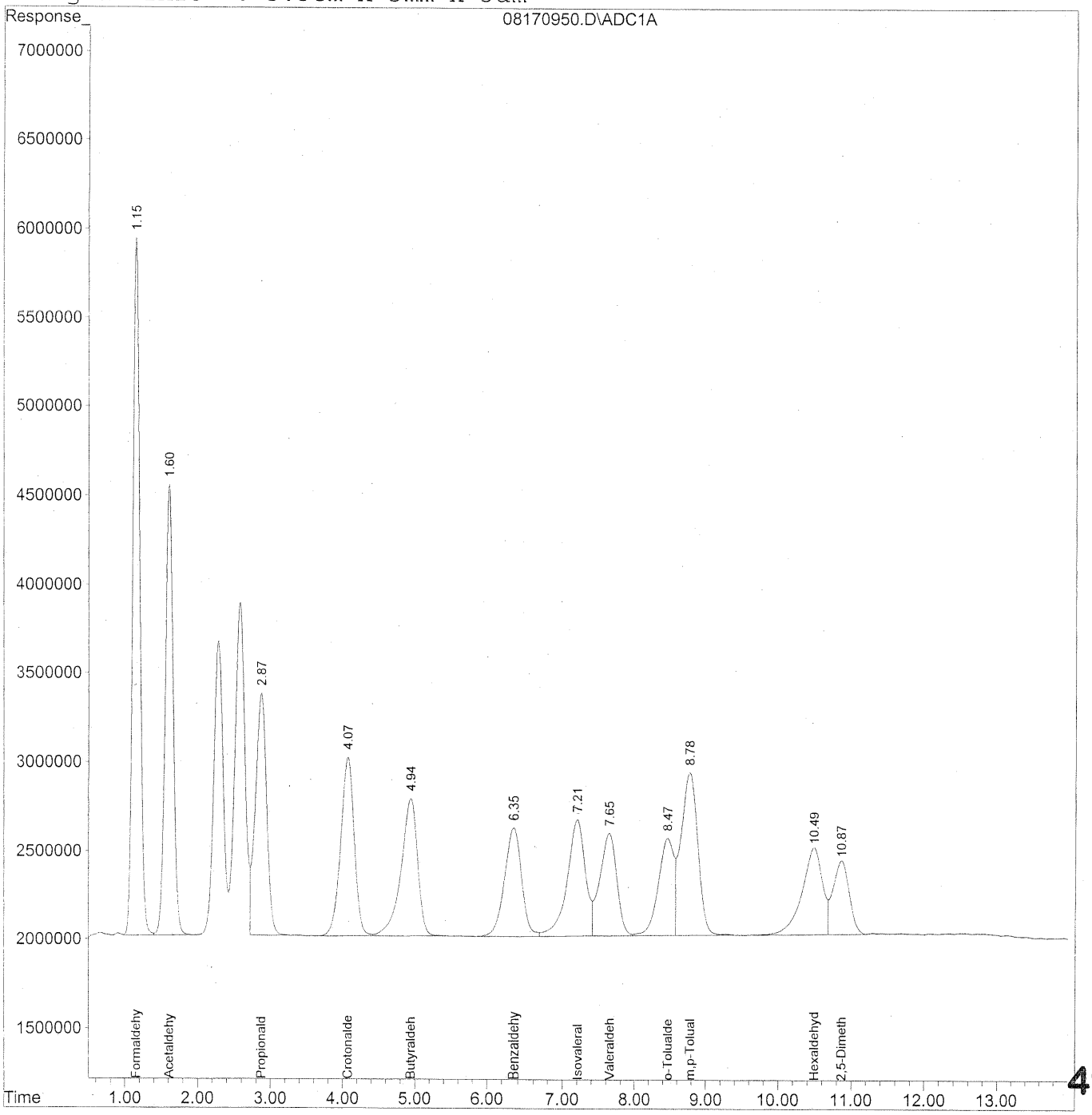
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170950.D Vial: 49
Acq On : 18 Aug 2009 3:07 am Operator: HC
Sample : CCV 1500ng/ml S21-08170901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 18 11:28 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 17 18:29:01 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



494

Quantitation Report (Not Reviewed)

QT HC 8/22/09

Data File : J:\LC01\DATA\TO11\2009_08\17\08170950.D Vial: 49
Acq On : 18 Aug 2009 3:07 am Operator: HC
Sample : CCV 1500ng/ml S21-08170901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 18 11:28 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 17 18:29:01 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

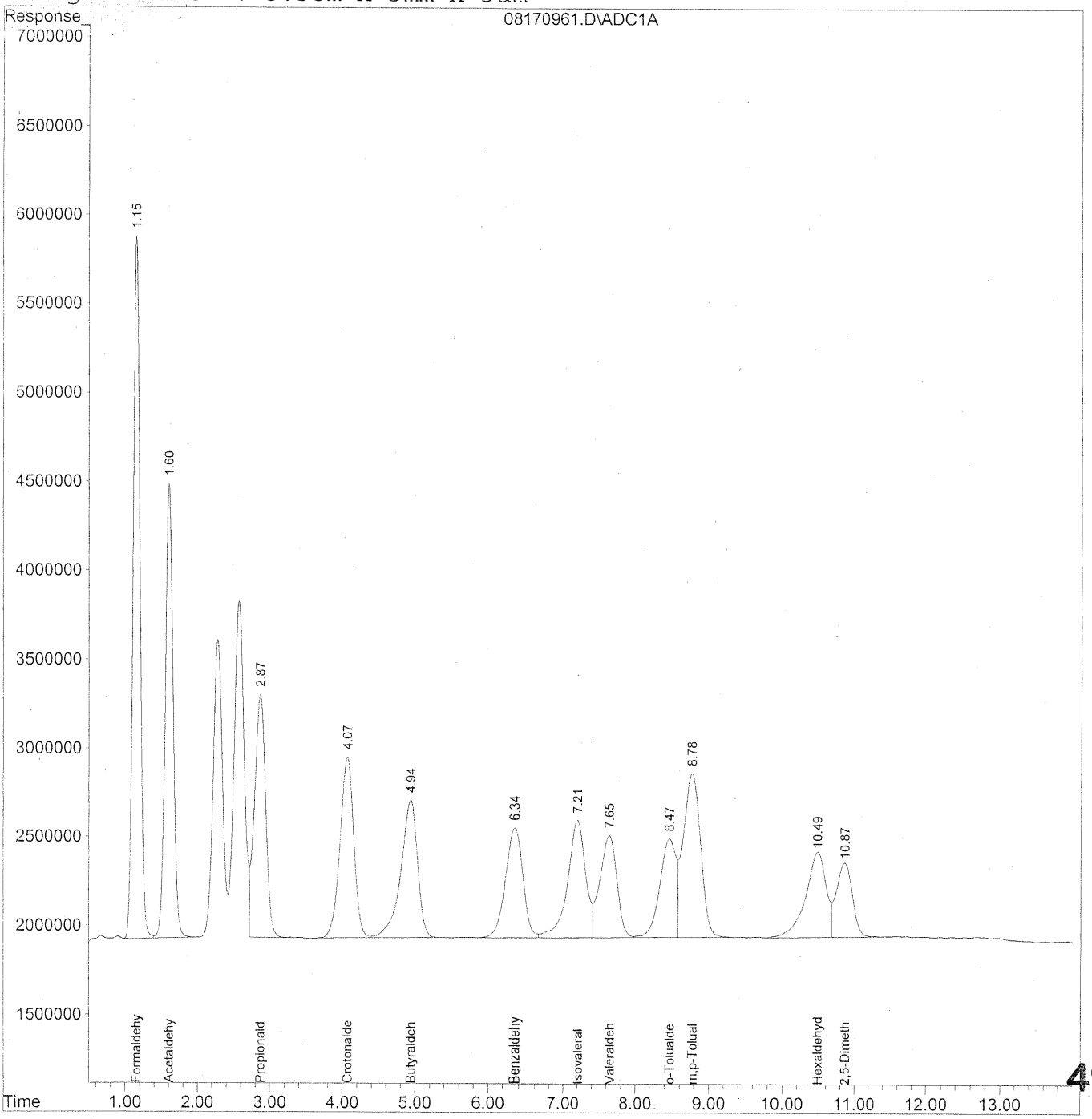
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.15	259780594	1415.070	ng/ml
2)	Acetaldehyde	1.60	194702902	1388.518	ng/ml
3)	Propionaldehyde	2.87	147496656	1382.411	ng/ml
4)	Crotonaldehyde	4.07	132496871	1360.126	ng/ml
5)	Butyraldehyde	4.94	124066617	1404.483	ng/ml
6)	Benzaldehyde	6.35	92617302	1406.077	ng/ml
7)	Isovaleraldehyde	7.21	115543030	1476.570	ng/ml
8)	Valeraldehyde	7.65	95237364	1295.658	ng/ml
9)	o-Tolualdehyde	8.47	83922090	1438.983	ng/ml
10)	m,p-Tolualdehyde	8.78	150902470	2794.728	ng/ml
11)	Hexaldehyde	10.49	95986938	1425.328	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.87	64584498	1317.691	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170961.D Vial: 59
Acq On : 18 Aug 2009 5:52 am Operator: HC
Sample : CCV 1500ng/ml S21-08170901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 18 8:06 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 17 18:29:01 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



496

Quantitation Report (Not Reviewed)

QT

Data File : J:\LC01\DATA\TO11\2009_08\17\08170961.D Vial: 59
Acq On : 18 Aug 2009 5:52 am Operator: HC
Sample : CCV 1500ng/ml S21-08170901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 18 8:06 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 17 18:29:01 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

HC
8/22/09

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

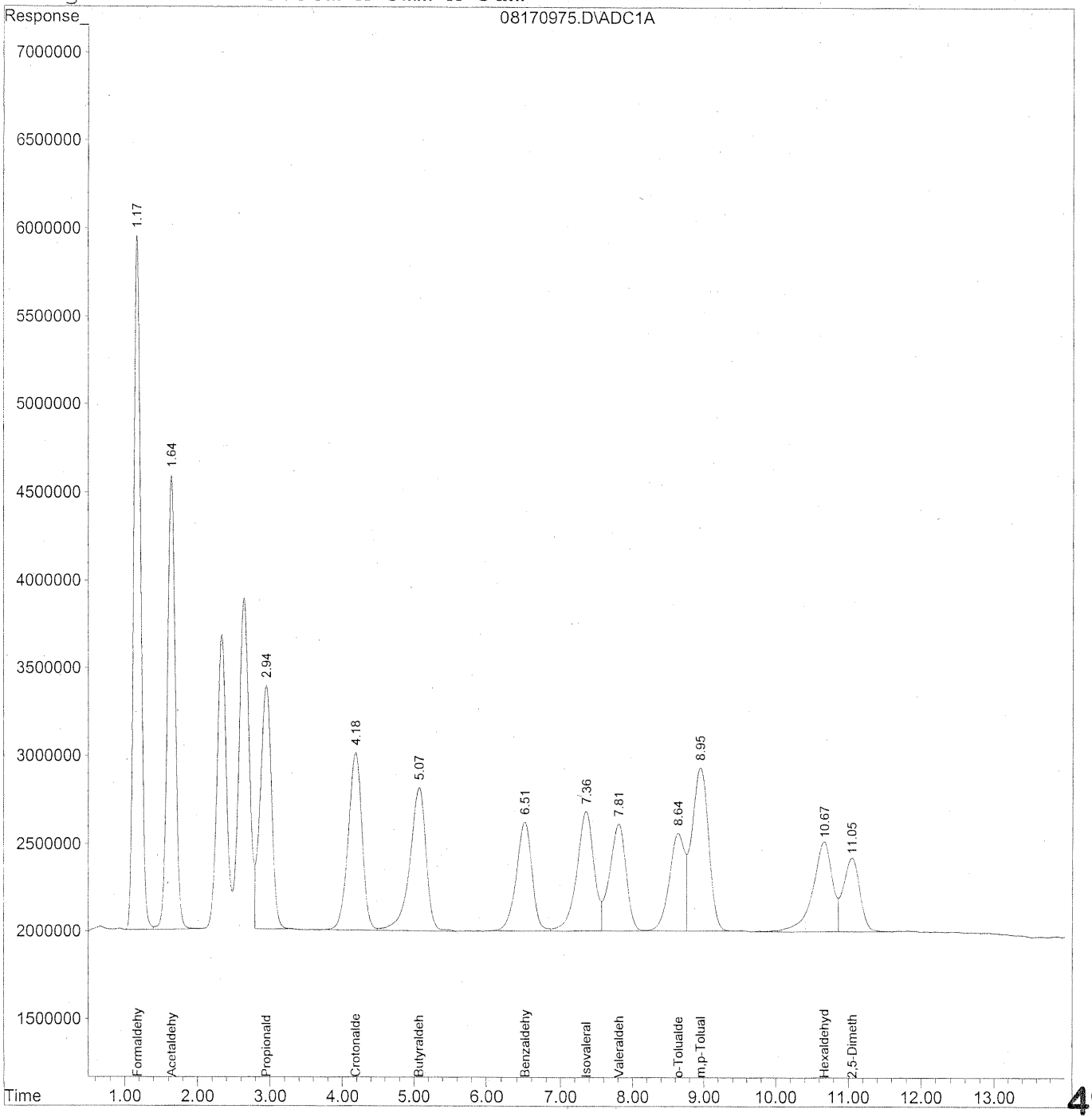
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.15	261296664	1423.329	ng/ml
2)	Acetaldehyde	1.60	196847529	1403.813	ng/ml
3)	Propionaldehyde	2.87	147953821	1386.696	ng/ml
4)	Crotonaldehyde	4.07	133538675	1370.820	ng/ml
5)	Butyraldehyde	4.94	125041065	1415.515	ng/ml
6)	Benzaldehyde	6.35	92503323	1404.346	ng/ml
7)	Isovaleraldehyde	7.21	116010499	1482.544	ng/ml
8)	Valeraldehyde	7.65	93732904	1275.191	ng/ml
9)	o-Tolualdehyde	8.47	84024696	1440.742	ng/ml
10)	m,p-Tolualdehyde	8.78	151692317	2809.356	ng/ml
11)	Hexaldehyde	10.49	95193329	1413.543	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.87	65738993	1341.246	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08170975.D Vial: 73
Acq On : 18 Aug 2009 9:23 am Operator: HC
Sample : CCV 1500ng/ml S21-08170902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:02 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



498

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170975.D Vial: 73
Acq On : 18 Aug 2009 9:23 am Operator: HC
Sample : CCV 1500ng/ml S21-08170902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:02 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

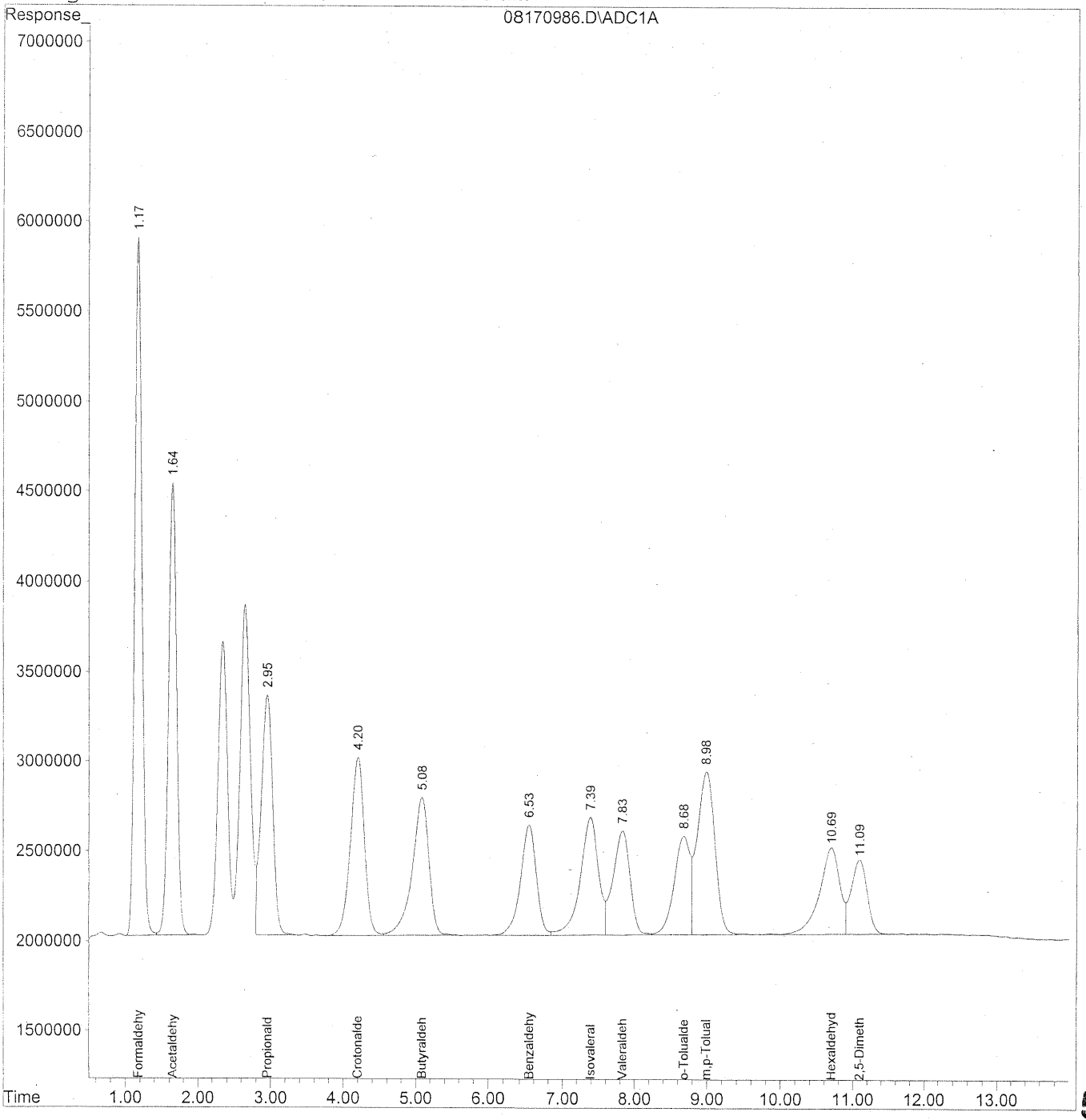
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	265760309	1447.643	ng/ml
2)	Acetaldehyde	1.64	201701487	1438.428	ng/ml
3)	Propionaldehyde	2.94	150378331	1409.420	ng/ml
4)	Crotonaldehyde	4.18	134736271	1383.114	ng/ml
5)	Butyraldehyde	5.07	128699146	1456.925	ng/ml
6)	Benzaldehyde	6.51	92796870	1408.803	ng/ml
7)	Isovaleraldehyde	7.36	113737485	1453.496	ng/ml
8)	Valeraldehyde	7.81	97709233	1329.287	ng/ml
9)	o-Tolualdehyde	8.64	83845925	1437.677	ng/ml
10)	m,p-Tolualdehyde	8.95	152620527	2826.547	ng/ml
11)	Hexaldehyde	10.67	98261245	1459.099	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.05	67018720	1367.355	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Data File : J:\LC01\DATA\TO11\2009_08\17\08170986.D Vial: 83
Acq On : 18 Aug 2009 12:08 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:02 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



500

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08170986.D Vial: 83
Acq On : 18 Aug 2009 12:08 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:02 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

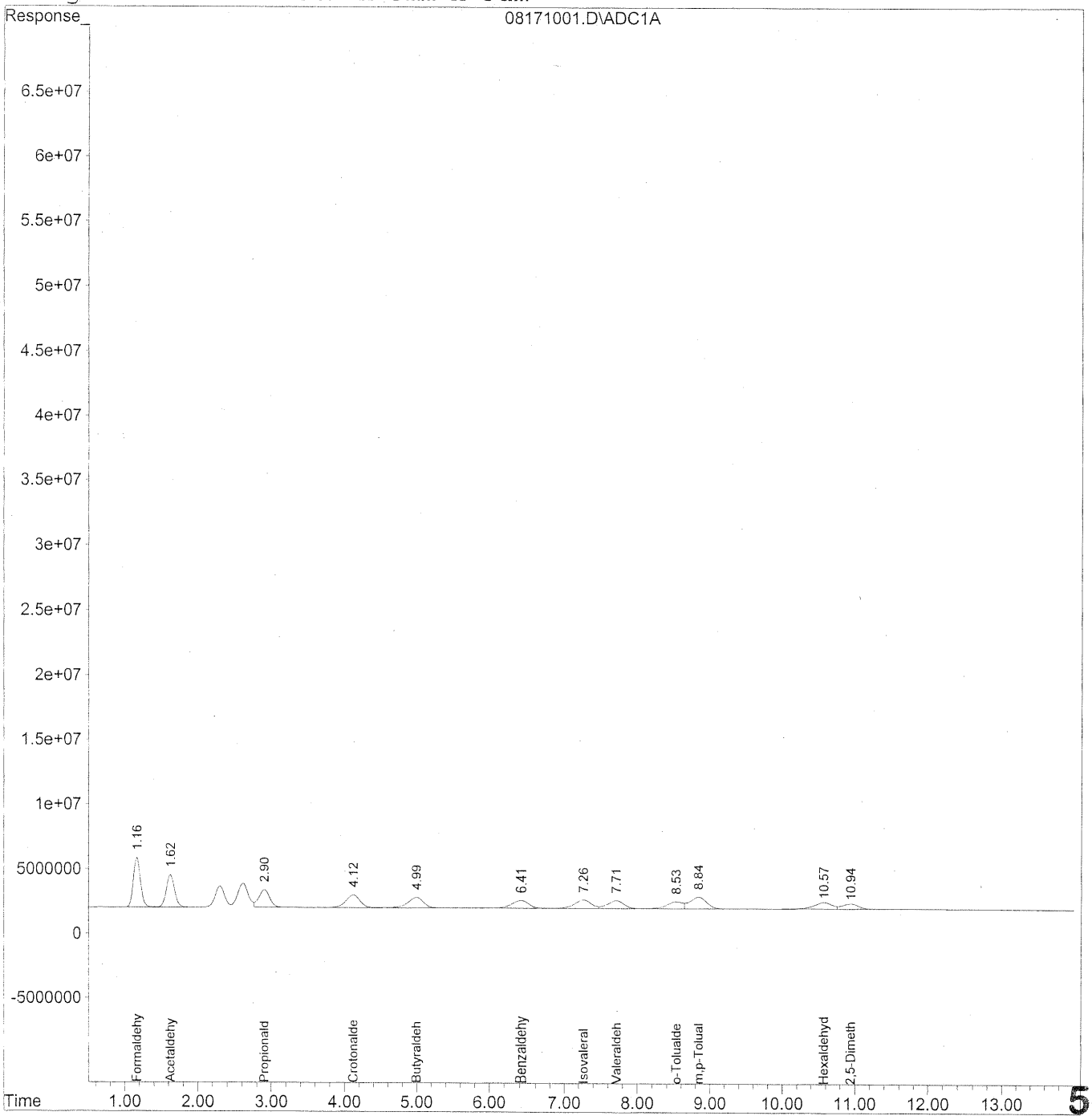
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	258583568	1408.550	ng/ml
2)	Acetaldehyde	1.64	195082035	1391.222	ng/ml
3)	Propionaldehyde	2.95	146615880	1374.156	ng/ml
4)	Crotonaldehyde	4.20	133152538	1366.856	ng/ml
5)	Butyraldehyde	5.08	124002584	1403.759	ng/ml
6)	Benzaldehyde	6.53	91636990	1391.194	ng/ml
7)	Isovaleraldehyde	7.39	113976379	1456.549	ng/ml
8)	Valeraldehyde	7.83	95604953	1300.659	ng/ml
9)	o-Tolualdehyde	8.68	82891702	1421.315	ng/ml
10)	m,p-Tolualdehyde	8.99	150699951	2790.978	ng/ml
11)	Hexaldehyde	10.70	93760303	1392.264	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.09	64640922	1318.842	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

Data File : J:\LC01\DATA\TO11\2009_08\17\08171001.D Vial: 15
Acq On : 18 Aug 2009 3:54 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 24 17:38 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 17 18:29:01 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



502

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171001.D Vial: 15
Acq On : 18 Aug 2009 3:54 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 24 17:38 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 17 18:29:01 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

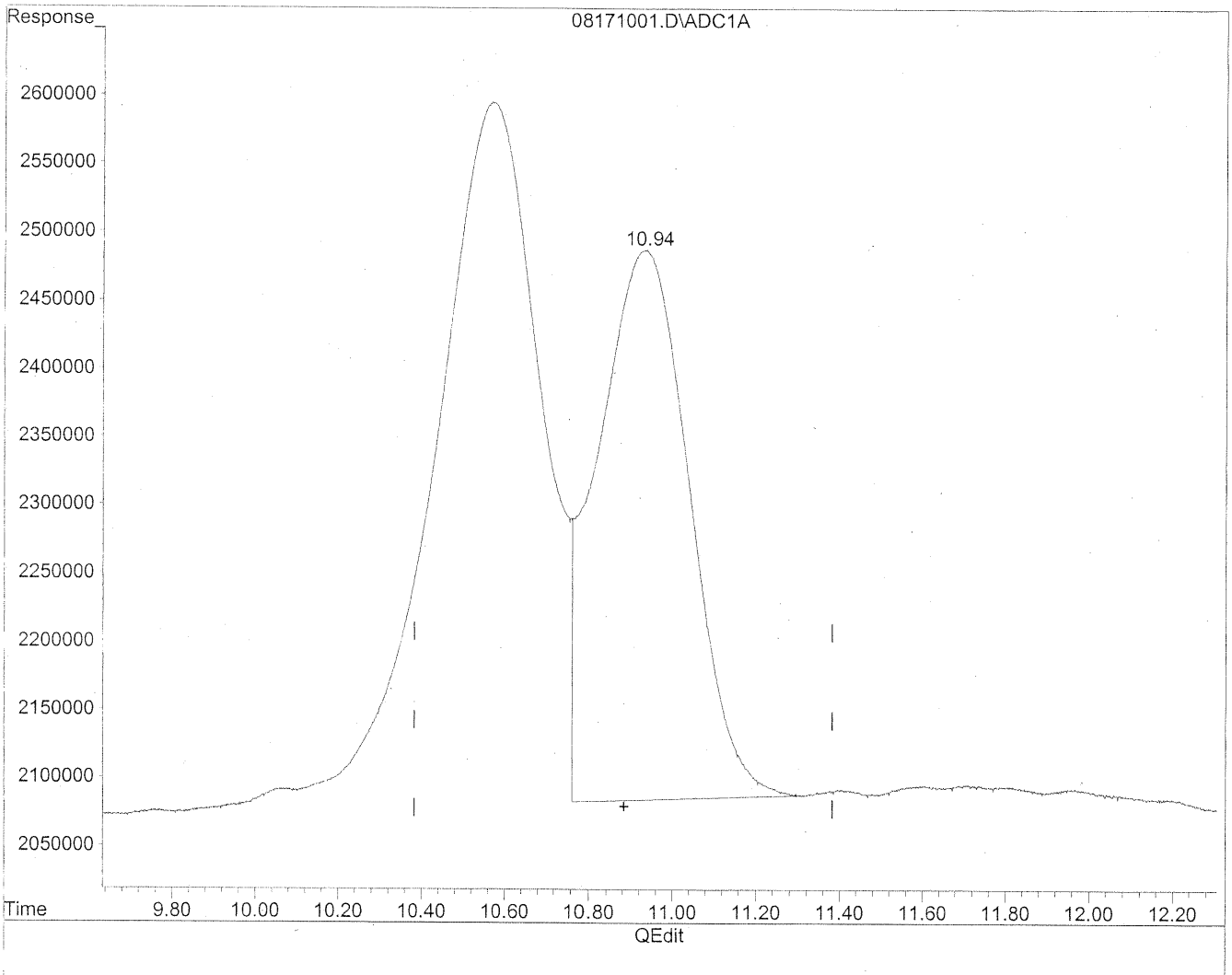
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	254492381	1386.265	ng/ml
2)	Acetaldehyde	1.62	192731672	1374.461	ng/ml
3)	Propionaldehyde	2.90	143194862	1342.093	ng/ml
4)	Crotonaldehyde	4.12	130784235	1342.545	ng/ml
5)	Butyraldehyde	4.99	123143394	1394.032	ng/ml
6)	Benzaldehyde	6.41	91322491	1386.419	ng/ml
7)	Isovaleraldehyde	7.26	108172704	1382.381	ng/ml
8)	Valeraldehyde	7.71	96087459	1307.223	ng/ml
9)	o-Tolualdehyde	8.53	81385268	1395.485	ng/ml
10)	m,p-Tolualdehyde	8.84	147203032	2726.214	ng/ml
11)	Hexaldehyde	10.57	92228372	1369.516	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.94	62912781	1283.584	ng/mlm

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171001.D Vial: 15
Acq On : 18 Aug 2009 3:54 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 18 16:10 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.93min 1260.162ng/ml

response 61764791

HC 8/24/09
RC no before
KP 8/25/09

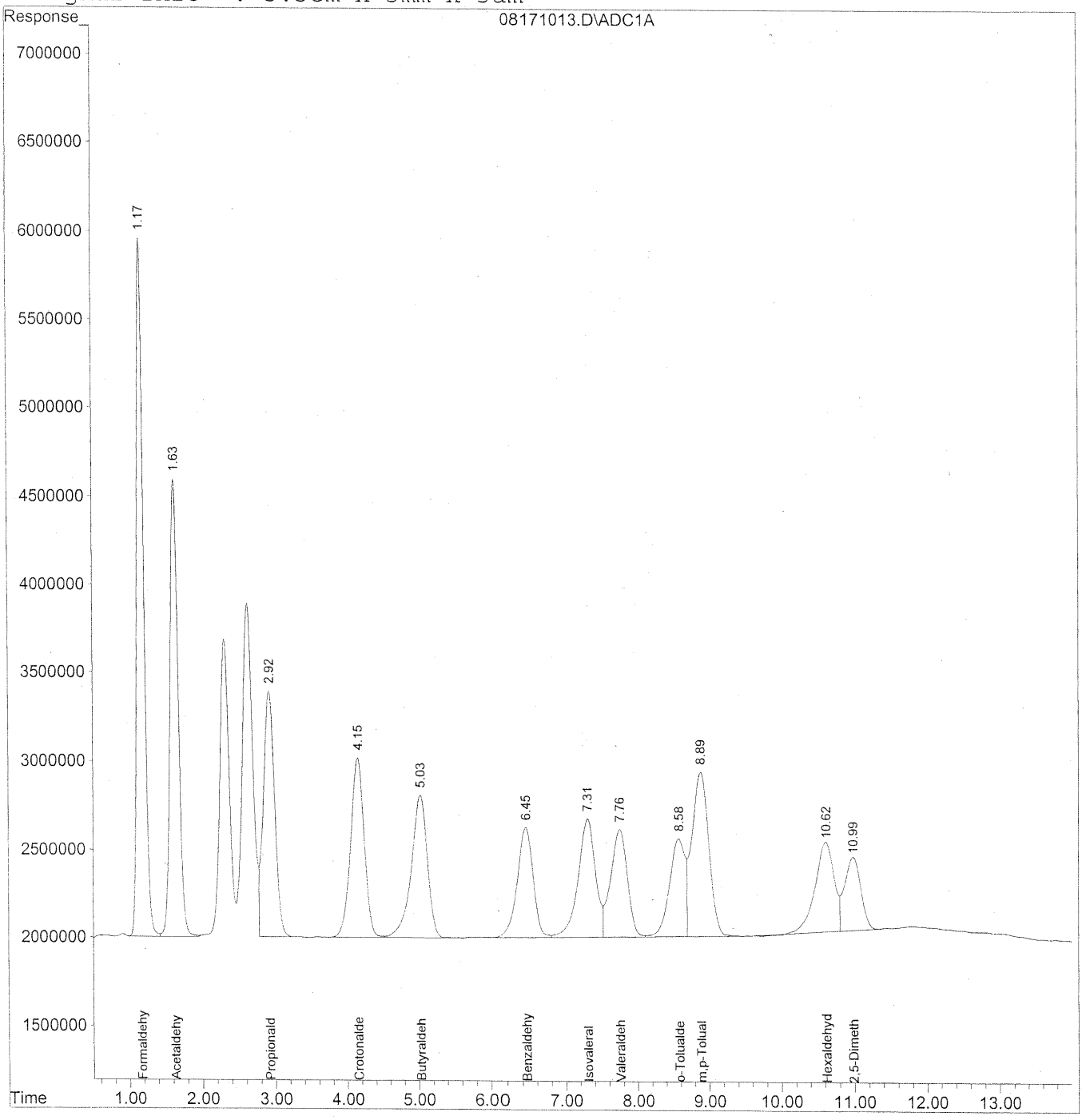
504

(+) = Expected Retention Time
08171001.D TO110709.M Mon Aug 24 17:38:22 2009

Data File : J:\LC01\DATA\TO11\2009_08\17\08171013.D Vial: 26
Acq On : 18 Aug 2009 6:54 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 8:59 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 18 17:12:05 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171013.D Vial: 26
Acq On : 18 Aug 2009 6:54 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 8:59 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 18 17:12:05 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	265986167	1448.873	ng/ml
2)	Acetaldehyde	1.63	202718555	1445.682	ng/ml
3)	Propionaldehyde	2.92	150704452	1412.476	ng/ml
4)	Crotonaldehyde	4.15	135136187	1387.219	ng/ml
5)	Butyraldehyde	5.03	126263286	1429.351	ng/ml
6)	Benzaldehyde	6.45	93091270	1413.272	ng/ml
7)	Isovaleraldehyde	7.31	113795597	1454.238	ng/ml
8)	Valeraldehyde	7.76	98983708	1346.625	ng/ml
9)	o-Tolualdehyde	8.58	85390518	1464.162	ng/ml
10)	m,p-Tolualdehyde	8.89	154221085	2856.189	ng/ml
11)	Hexaldehyde	10.61	93290039	1385.281	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.98	64366262	1313.238	ng/ml

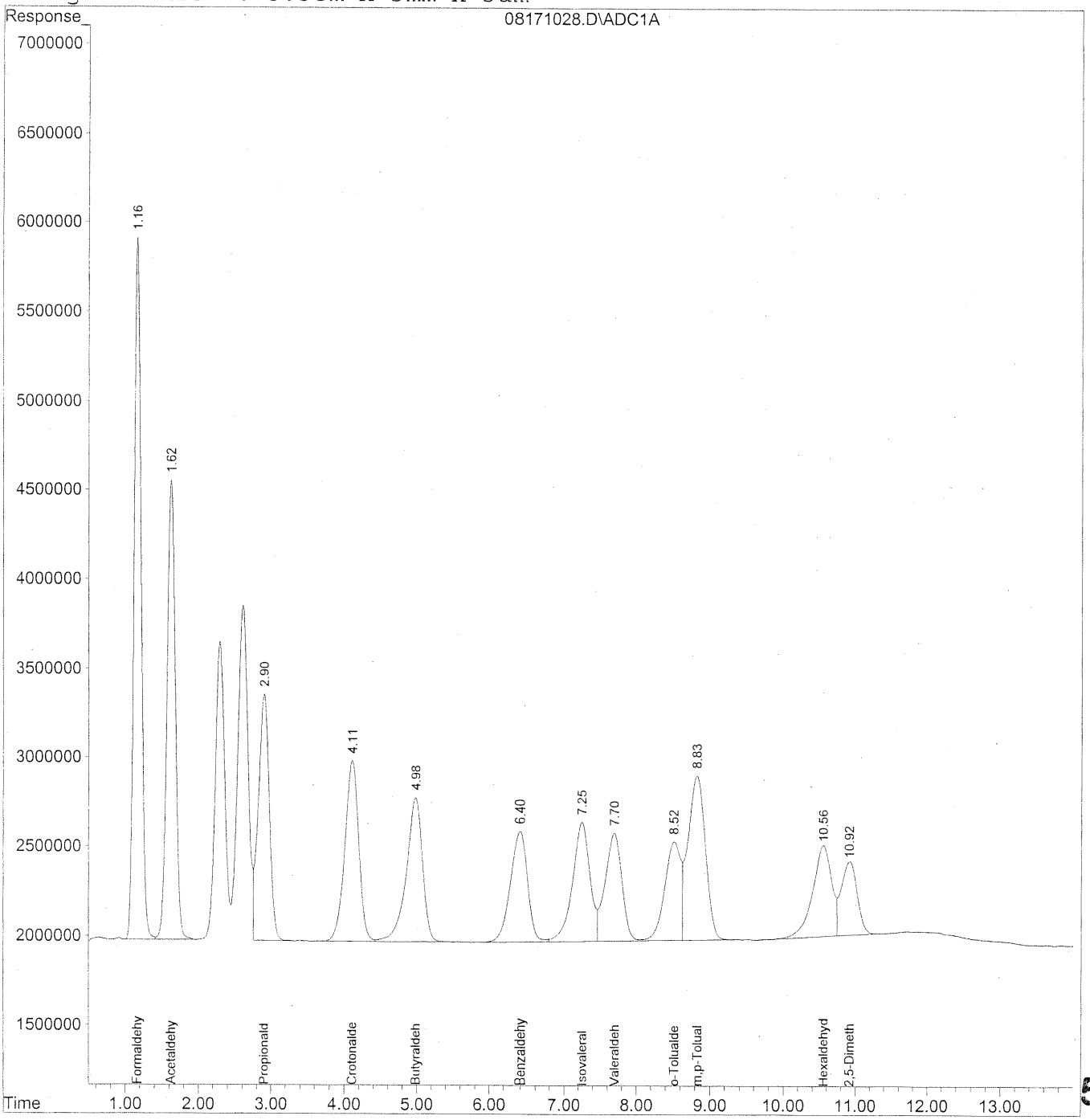
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171028.D Vial: 41
Acq On : 18 Aug 2009 10:40 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 8:58 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 18 17:12:05 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



507

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171028.D Vial: 41
Acq On : 18 Aug 2009 10:40 pm Operator: HC
Sample : CCV 1500ng/ml S21-08170902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 8:58 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 18 17:12:05 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

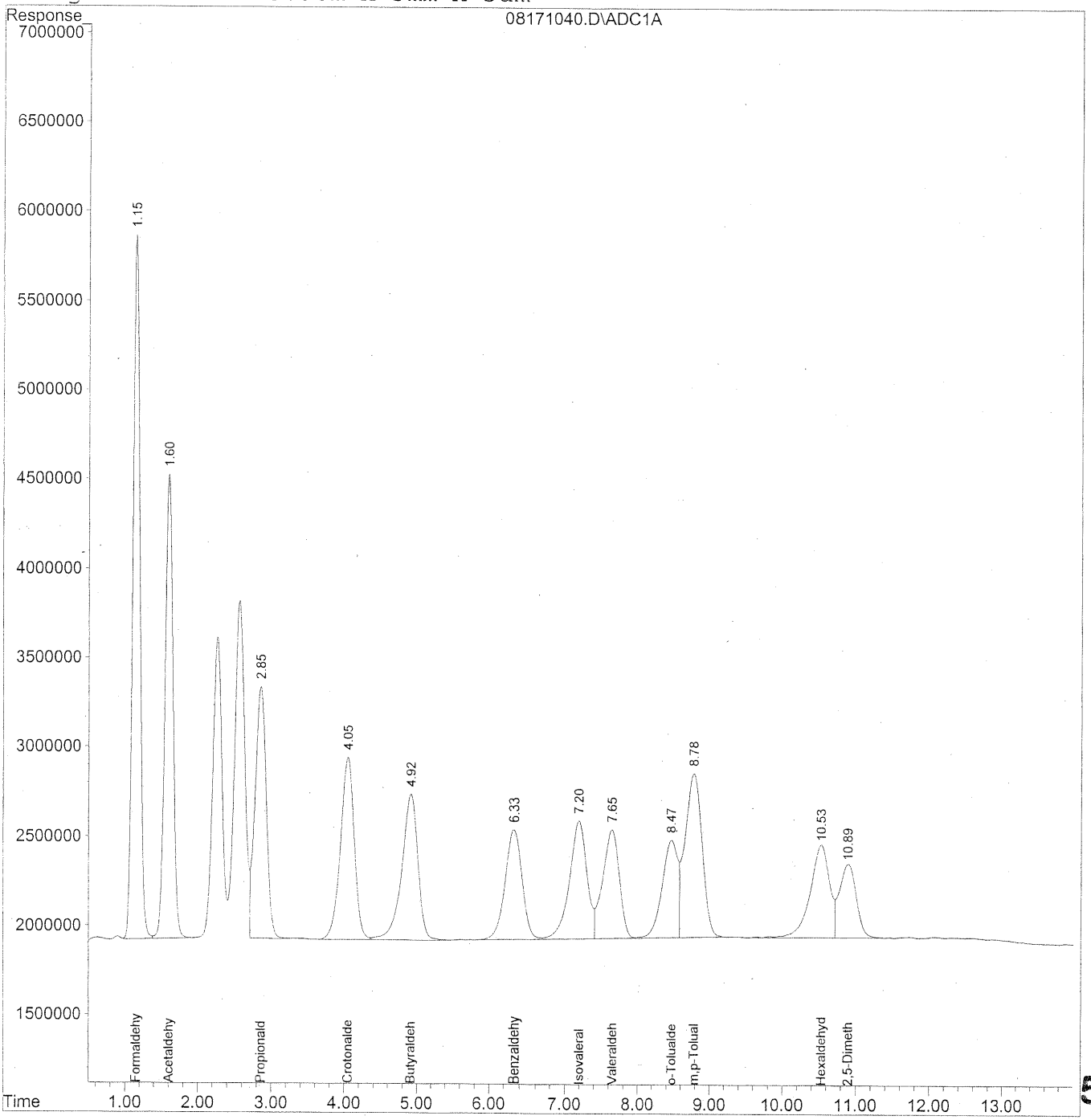
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	261871579	1426.460	ng/ml
2)	Acetaldehyde	1.62	198710587	1417.099	ng/ml
3)	Propionaldehyde	2.90	148575631	1392.524	ng/ml
4)	Crotonaldehyde	4.11	133760810	1373.100	ng/ml
5)	Butyraldehyde	4.98	125729792	1423.311	ng/ml
6)	Benzaldehyde	6.40	92597408	1405.775	ng/ml
7)	Isovaleraldehyde	7.25	112286059	1434.947	ng/ml
8)	Valeraldehyde	7.70	96968687	1319.212	ng/ml
9)	o-Tolualdehyde	8.52	83682539	1434.876	ng/ml
10)	m,p-Tolualdehyde	8.83	150779313	2792.447	ng/ml
11)	Hexaldehyde	10.55	94971176	1410.245	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.92	63580105	1297.199	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171040.D Vial: 52
Acq On : 19 Aug 2009 1:40 am Operator: HC
Sample : CCV 1500ng/ml S21-08170902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 8:56 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 18 17:12:05 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



509

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171040.D Vial: 52
Acq On : 19 Aug 2009 1:40 am Operator: HC
Sample : CCV 1500ng/ml S21-08170902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 8:56 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 18 17:12:05 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

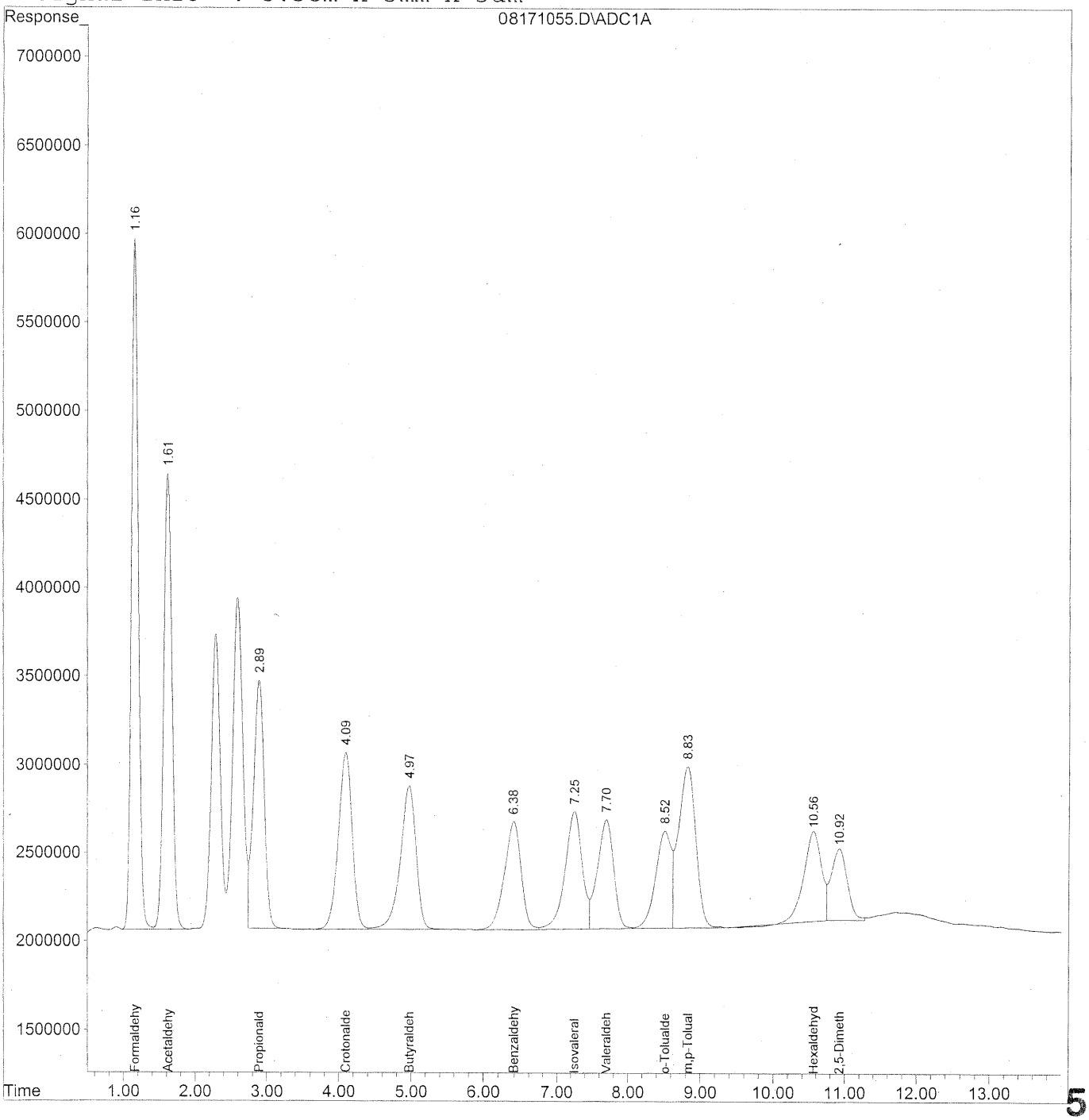
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.15	260757216	1420.390	ng/ml
2)	Acetaldehyde	1.60	196681817	1402.631	ng/ml
3)	Propionaldehyde	2.85	148836235	1394.966	ng/ml
4)	Crotonaldehyde	4.05	134361650	1379.268	ng/ml
5)	Butyraldehyde	4.92	128030206	1449.353	ng/ml
6)	Benzaldehyde	6.33	91638273	1391.213	ng/ml
7)	Isovaleraldehyde	7.20	110959211	1417.991	ng/ml
8)	Valeraldehyde	7.65	96909923	1318.413	ng/ml
9)	o-Tolualdehyde	8.47	83770663	1436.387	ng/ml
10)	m,p-Tolualdehyde	8.78	150515020	2787.553	ng/ml
11)	Hexaldehyde	10.53	97556071	1448.628	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.90	65975397	1346.069	ng/ml

Data File : J:\LC01\DATA\TO11\2009_08\17\08171055.D Vial: 67
Acq On : 19 Aug 2009 5:26 am Operator: HC
Sample : CCV 1500ng/ml S21-08180901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:05 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 18 17:12:05 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171055.D Vial: 67
Acq On : 19 Aug 2009 5:26 am Operator: HC
Sample : CCV 1500ng/ml S21-08180901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 9:05 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 18 17:12:05 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

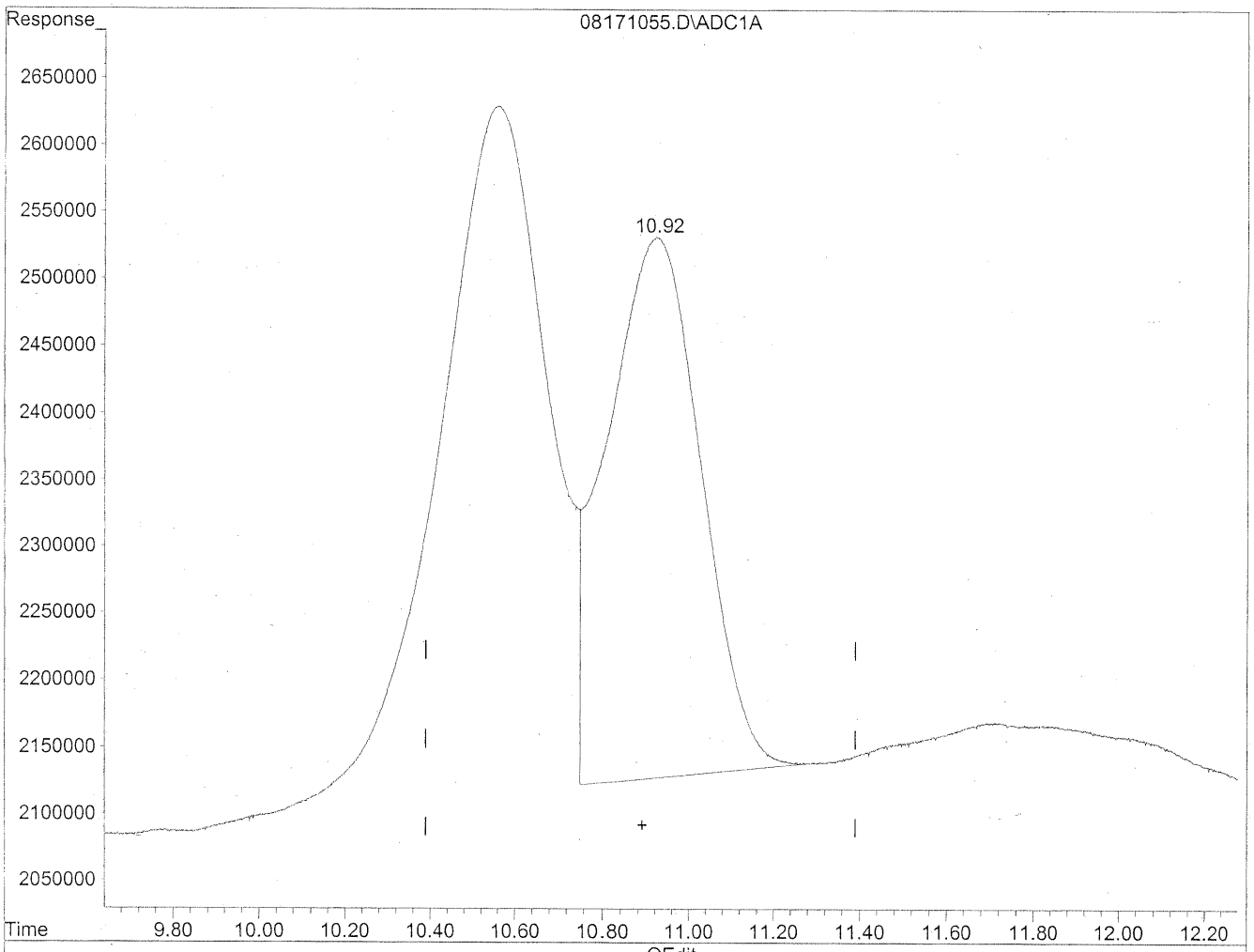
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	261118258	1422.357	ng/ml
2)	Acetaldehyde	1.62	197102998	1405.634	ng/ml
3)	Propionaldehyde	2.89	148877470	1395.353	ng/ml
4)	Crotonaldehyde	4.09	133086274	1366.176	ng/ml
5)	Butyraldehyde	4.97	124804286	1412.834	ng/ml
6)	Benzaldehyde	6.39	92714008	1407.545	ng/ml
7)	Isovaleraldehyde	7.25	112459019	1437.158	ng/ml
8)	Valeraldehyde	7.70	98515624	1340.257	ng/ml
9)	o-Tolualdehyde	8.52	83693115	1435.057	ng/ml
10)	m,p-Tolualdehyde	8.83	152140194	2817.651	ng/ml
11)	Hexaldehyde	10.56	89509172	1329.138	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.92	64421120	1314.358	ng/mlm

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171055.D Vial: 67
Acq On : 19 Aug 2009 5:26 am Operator: HC
Sample : CCV 1500ng/ml S21-08180901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 8:55 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.92min 1253.740ng/ml

response 61450053

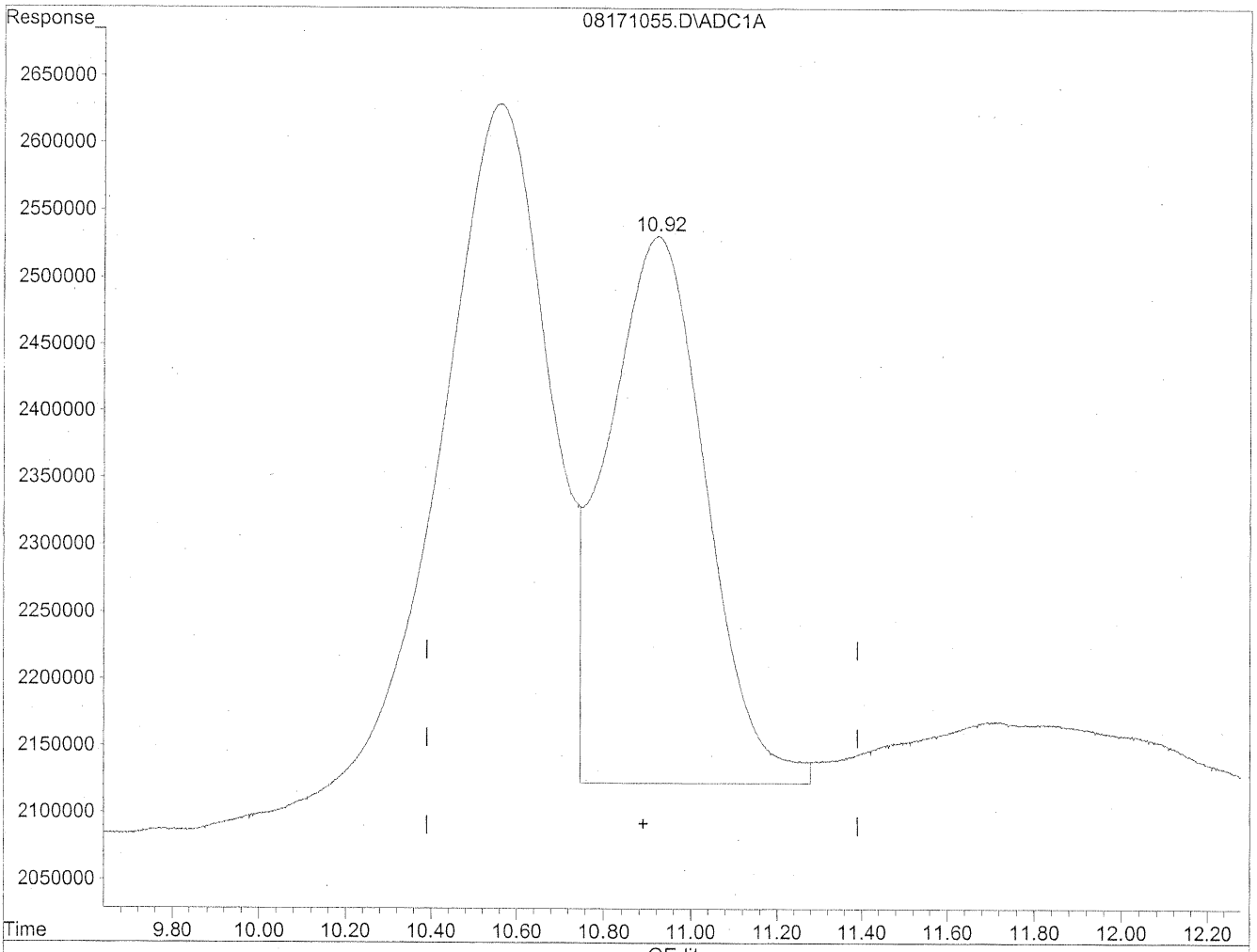
(+) = Expected Retention Time

08171055.D TO110709.M Wed Aug 19 09:04:47 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\17\08171055.D Vial: 67
Acq On : 19 Aug 2009 5:26 am Operator: HC
Sample : CCV 1500ng/ml S21-08180901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 8:55 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 09:01:59 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.92min 1314.358ng/ml m

response 64421120

8/19/09

8/23/09

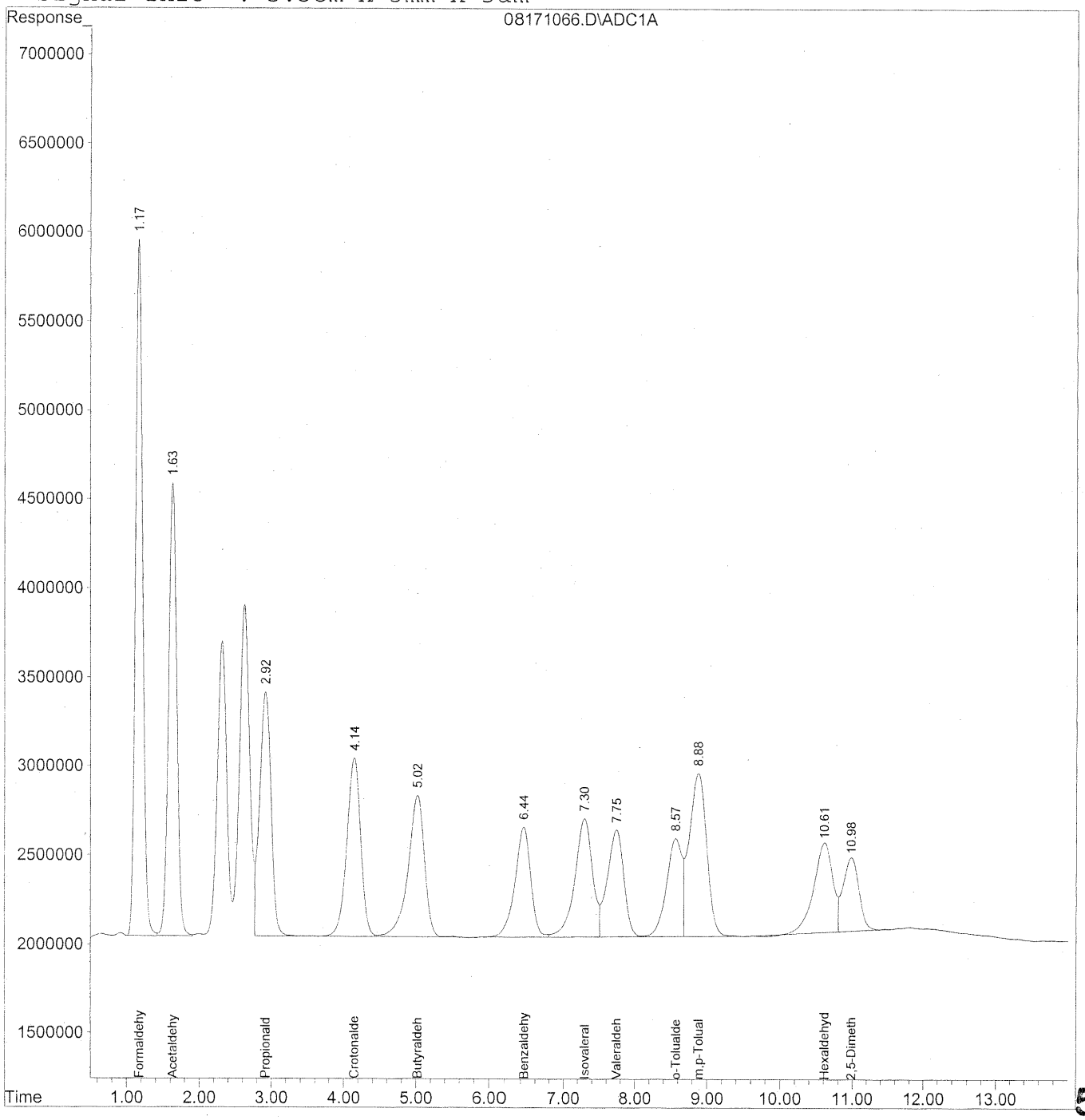
514

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_08\17\08171066.D Vial: 77
Acq On : 19 Aug 2009 8:11 am Operator: HC
Sample : CCV 1500ng/ml S21-08180901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 8:54 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 18 17:12:05 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



515

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171066.D Vial: 77
Acq On : 19 Aug 2009 8:11 am Operator: HC
Sample : CCV 1500ng/ml S21-08180901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 8:54 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 18 17:12:05 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

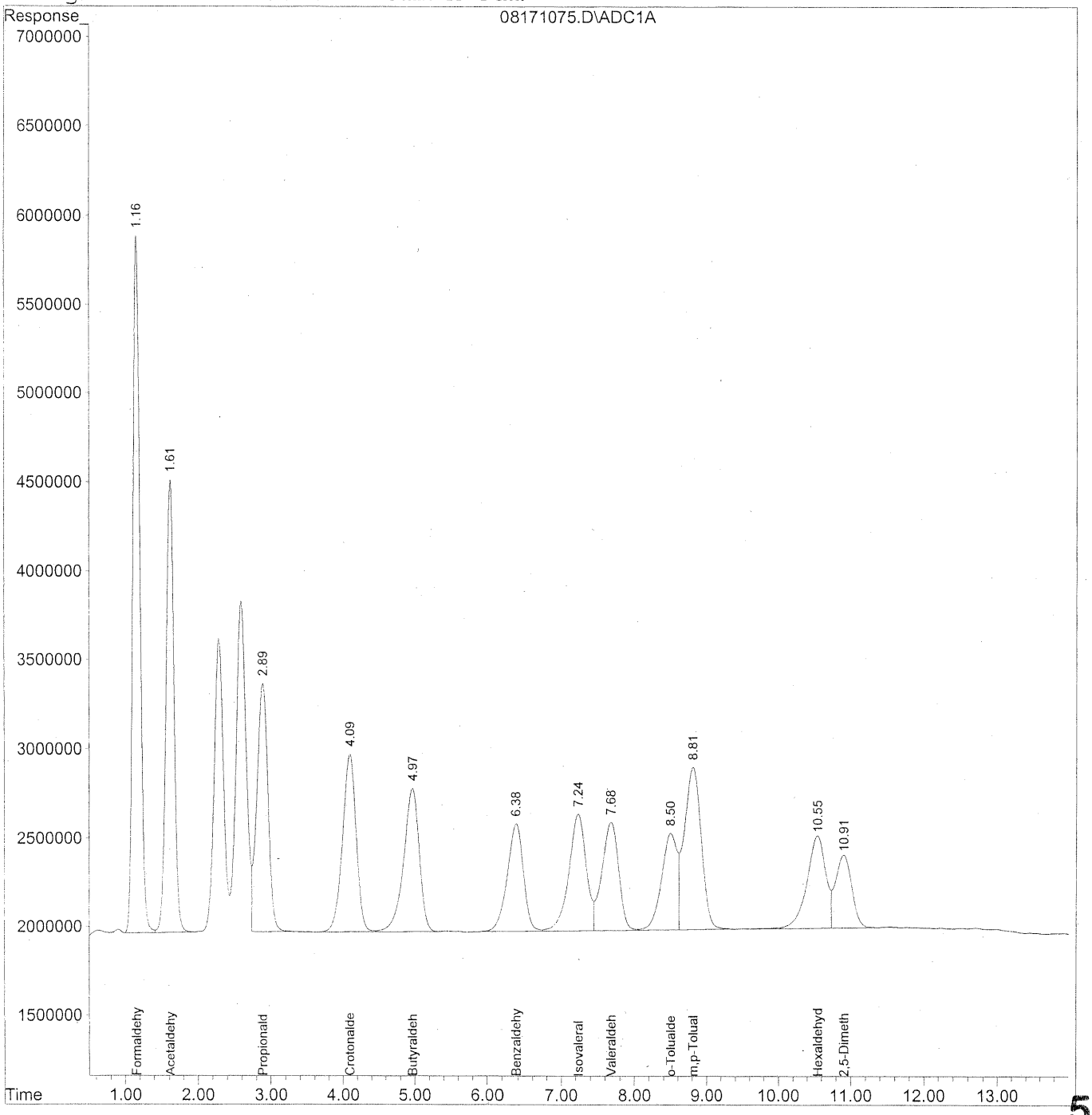
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	260925407	1421.306	ng/ml
2)	Acetaldehyde	1.63	197643204	1409.487	ng/ml
3)	Propionaldehyde	2.92	149366042	1399.932	ng/ml
4)	Crotonaldehyde	4.14	133349423	1368.877	ng/ml
5)	Butyraldehyde	5.02	124000936	1403.740	ng/ml
6)	Benzaldehyde	6.44	91469310	1388.648	ng/ml
7)	Isovaleraldehyde	7.30	113036582	1444.539	ng/ml
8)	Valeraldehyde	7.74	96202864	1308.793	ng/ml
9)	o-Tolualdehyde	8.57	83394674	1429.940	ng/ml
10)	m,p-Tolualdehyde	8.88	152093275	2816.782	ng/ml
11)	Hexaldehyde	10.61	91558527	1359.570	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.98	64026808	1306.313	ng/ml

Data File : J:\LC01\DATA\TO11\2009_08\17\08171075.D Vial: 85
Acq On : 19 Aug 2009 10:26 am Operator: HC
Sample : CCV 1500ng/ml S21-08180901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 10:45 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Data File : J:\LC01\DATA\TO11\2009_08\17\08171075.D Vial: 85
Acq On : 19 Aug 2009 10:26 am Operator: HC
Sample : CCV 1500ng/ml S21-08180901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 10:45 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

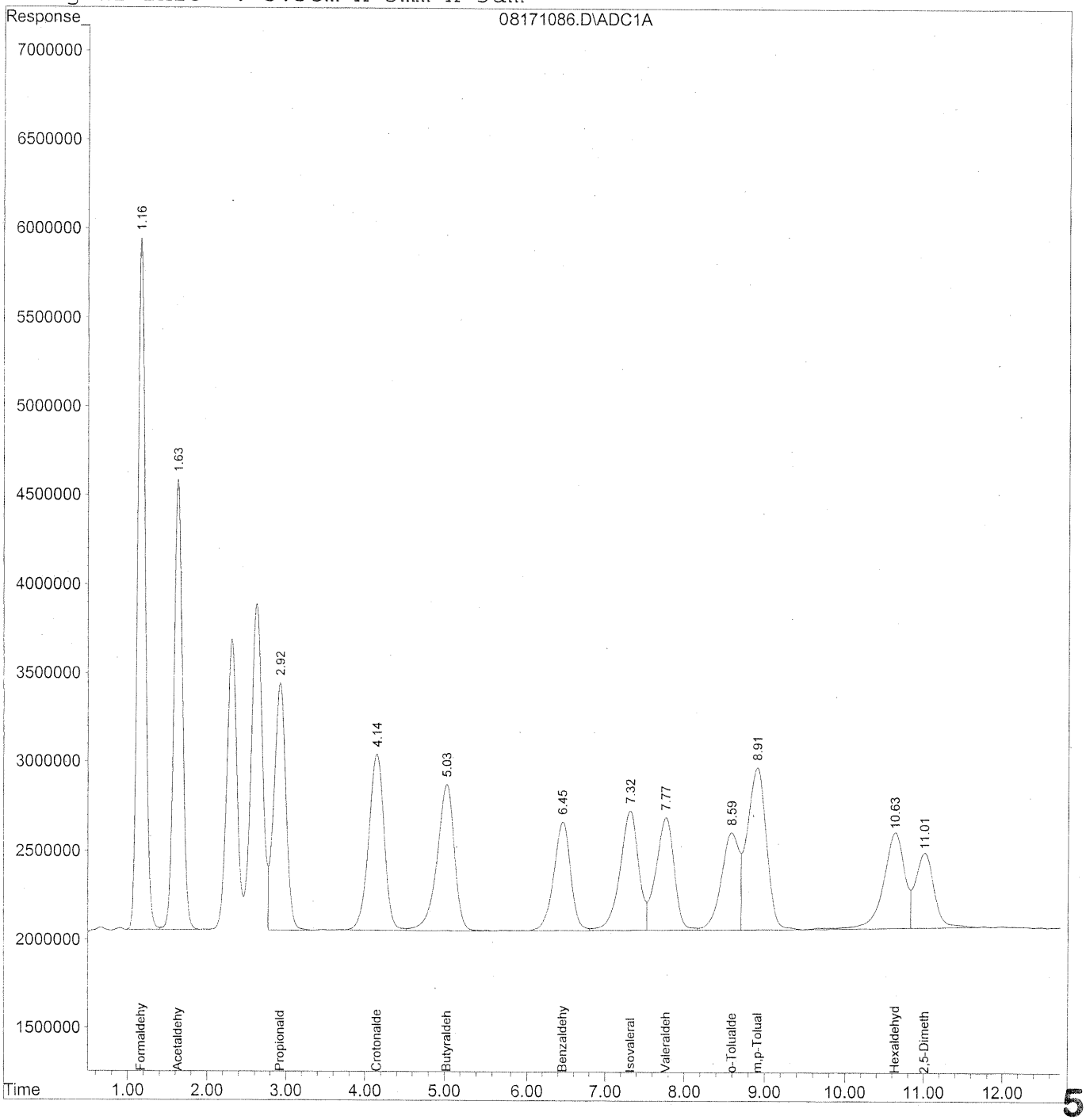
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	260984531	1421.628	ng/ml
2)	Acetaldehyde	1.61	195283064	1392.656	ng/ml
3)	Propionaldehyde	2.89	148529239	1392.089	ng/ml
4)	Crotonaldehyde	4.09	132676484	1361.969	ng/ml
5)	Butyraldehyde	4.97	122913452	1391.429	ng/ml
6)	Benzaldehyde	6.38	89995023	1366.266	ng/ml
7)	Isovaleraldehyde	7.24	109529955	1399.726	ng/ml
8)	Valeraldehyde	7.69	97192463	1322.256	ng/ml
9)	o-Tolualdehyde	8.50	82265870	1410.584	ng/ml
10)	m,p-Tolualdehyde	8.81	151072803	2797.883	ng/ml
11)	Hexaldehyde	10.54	96204294	1428.555	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.91	64218471	1310.223	ng/ml

Data File : J:\LC01\DATA\TO11\2009_08\17\08171086.D Vial: 11
Acq On : 19 Aug 2009 1:12 pm Operator: HC
Sample : CCV 1500ng/ml S21-08180901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 13:24 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\17\08171086.D Vial: 11
Acq On : 19 Aug 2009 1:12 pm Operator: HC
Sample : CCV 1500ng/ml S21-08180901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 19 13:24 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Wed Aug 19 10:45:48 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	259423883	1413.127	ng/ml
2)	Acetaldehyde	1.63	194867352	1389.691	ng/ml
3)	Propionaldehyde	2.92	147991378	1387.048	ng/ml
4)	Crotonaldehyde	4.15	132045693	1355.494	ng/ml
5)	Butyraldehyde	5.03	125141421	1416.651	ng/ml
6)	Benzaldehyde	6.46	91802294	1393.704	ng/ml
7)	Isovaleraldehyde	7.31	110206985	1408.378	ng/ml
8)	Valeraldehyde	7.77	100536028	1367.744	ng/ml
9)	o-Tolualdehyde	8.59	83596705	1433.404	ng/ml
10)	m,p-Tolualdehyde	8.90	151854950	2812.368	ng/ml
11)	Hexaldehyde	10.64	101536821	1507.739	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.01	69602596	1420.073	ng/ml

(f)=RT Delta > 1/2 Window (m)=manual int.

RUN LOGS

Injection Log

Directory: j:\lc01\data\to11\2009_07\28

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	1	07280901.d	1.	Prime		28 Jul 109 12::3
2	2	07280902.d	1.	ACN blank Lot CY023		28 Jul 109 12::5
3	3	07280903.d	1.	50ng/ml TO11A Std S21-07270908		28 Jul 109 12::0
4	4	07280904.d	1.	50ng/ml TO11A Std S21-07270908		28 Jul 109 12::2
5	5	07280905.d	1.	50ng/ml TO11A Std S21-07270908		28 Jul 109 12::3
6	6	07280906.d	1.	100ng/ml TO11A Std S21-07270905		28 Jul 109 12::5
7	7	07280907.d	1.	100ng/ml TO11A Std S21-07270905		28 Jul 109 13::0
8	8	07280908.d	1.	100ng/ml TO11A Std S21-07270905		28 Jul 109 13::2
9	9	07280909.d	1.	500ng/ml TO11A Std S21-07270904		28 Jul 109 13::3
10	10	07280910.d	1.	500ng/ml TO11A Std S21-07270904		28 Jul 109 13::5
11	11	07280911.d	1.	500ng/ml TO11A Std S21-07270904		28 Jul 109 13::0
12	12	07280912.d	1.	1500ng/ml TO11A Std S21-07270903		28 Jul 109 13::2
13	13	07280913.d	1.	1500ng/ml TO11A Std S21-07270903		28 Jul 109 13::3
14	14	07280914.d	1.	1500ng/ml TO11A Std S21-07270903		28 Jul 109 13::5
15	15	07280915.d	1.	5000ng/ml TO11A Std S21-07270902		28 Jul 109 13::0
16	16	07280916.d	1.	5000ng/ml TO11A Std S21-07270902		28 Jul 109 13::2
17	17	07280917.d	1.	5000ng/ml TO11A Std S21-07270902		28 Jul 109 13::3
18	18	07280918.d	1.	10000ng/ml TO11A Std S21-07270901		28 Jul 109 13::5
19	19	07280919.d	1.	10000ng/ml TO11A Std S21-07270901		28 Jul 109 12::0
20	20	07280920.d	1.	10000ng/ml TO11A Std S21-07270901		28 Jul 109 12::2
21	21	07280921.d	1.	~1500ng/ml TO11A Std ICV S21-07270907		28 Jul 109 12::4

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Injection Log

Directory: j:\lc01\data\to11\2009_08\17

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	1	08170901.d	1.	PRIME		17 Aug 109 12::
2	2	08170902.d	1.	1500ng/ml TO11A std S21-08170901		17 Aug 109 12::
3	3	08170903.d	1.	ACN blank Lot CY023		17 Aug 109 12::
4	4	08170904.d	1.	P0902771-008 front 10x		17 Aug 109 12::
5	5	08170905.d	1.	P0902771-010 front 10x		17 Aug 109 12::
6	6	08170906.d	1.	P0902771-011 front 10x		17 Aug 109 12::
7	7	08170907.d	1.	P0902771-018 front 10x		17 Aug 109 12::
8	8	08170908.d	1.	P0902771-021 front 10x		17 Aug 109 12::
9	9	08170909.d	1.	P0902771-022 front 10x		17 Aug 109 12::
10	10	08170910.d	1.	P0902771-024 front 10x		17 Aug 109 12::
11	11	08170911.d	1.	CCV 1500ng/ml S21-08170901		17 Aug 109 12::
12	12	08170912.d	1.	ACN CY023 blk		17 Aug 109 12::
13	13	08170913.d	1.	MB front lot 6009/6097 1.0ml		17 Aug 109 12::
14	14	08170914.d	1.	MB back lot 6009/6097 1.0ml		17 Aug 109 12::
15	15	08170915.d	1.	P0902770-001 back 1.0ml		17 Aug 109 12::
16	16	08170916.d	1.	P0902770-002 back 1.0ml		17 Aug 109 12::
17	17	08170917.d	1.	P0902770-003 back 1.0ml		17 Aug 109 12::
18	18	08170918.d	1.	P0902770-004 back 1.0ml		17 Aug 109 12::
19	19	08170919.d	1.	P0902770-005 back 1.0ml		17 Aug 109 12::
20	20	08170920.d	1.	P0902770-006 back 1.0ml		17 Aug 109 12::
21	21	08170921.d	1.	P0902770-007 back 1.0ml		17 Aug 109 12::
22	22	08170922.d	1.	P0902770-008 back 1.0ml		17 Aug 109 12::
23	23	08170923.d	1.	P0902770-009 back 1.0ml		17 Aug 109 12::
24	24	08170924.d	1.	P0902770-010 back 1.0ml		17 Aug 109 12::
25	25	08170925.d	1.	CCV 1500ng/ml S21-08170901		17 Aug 109 12::
26	26	08170926.d	1.	P0902770-011 back 1.0ml		17 Aug 109 12::
27	26	08170927.d	1.	P0902770-011dup back 1.0ml		17 Aug 109 12::
28	27	08170928.d	1.	P0902770-012 back 1.0ml		17 Aug 109 12::
29	28	08170929.d	1.	P0902770-013 back 1.0ml		17 Aug 109 12::
30	29	08170930.d	1.	P0902772-001 back 1.0ml		17 Aug 109 13::
31	30	08170931.d	1.	P0902772-002 back 1.0ml		17 Aug 109 13::
32	31	08170932.d	1.	P0902772-003 back 1.0ml		17 Aug 109 13::
33	32	08170933.d	1.	P0902772-004 back 1.0ml		17 Aug 109 13::
34	33	08170934.d	1.	P0902772-005 back 1.0ml		17 Aug 109 13::
35	34	08170935.d	1.	P0902772-006 back 1.0ml		17 Aug 109 13::
36	35	08170936.d	1.	CCV 1500ng/ml S21-08170901		17 Aug 109 13::
37	36	08170937.d	1.	ACN blk lot CY023		17 Aug 109 13::
38	37	08170938.d	1.	MB front lot 6009/6097 1.0ml		18 Aug 109 13::
39	38	08170939.d	1.	MB back lot 6009/6097 1.0ml		18 Aug 109 13::
40	39	08170940.d	1.	P0902772-007 back 1.0ml		18 Aug 109 13::
41	40	08170941.d	1.	P0902772-008 back 1.0ml		18 Aug 109 13::
42	41	08170942.d	1.	P0902772-009 back1.0ml		18 Aug 109 12::
43	42	08170943.d	1.	P0902772-010 back 1.0ml		18 Aug 109 12::
44	43	08170944.d	1.	P0902772-011 back 1.0ml		18 Aug 109 12::
45	44	08170945.d	1.	P0902772-012 back 1.0ml		18 Aug 109 12::
46	45	08170946.d	1.	P0902786-001 back 1.0ml		18 Aug 109 12::
47	46	08170947.d	1.	P0902786-002 back 1.0ml		18 Aug 109 12::
48	47	08170948.d	1.	P0902786-003 back 1.0ml		18 Aug 109 12::
49	48	08170949.d	1.	P0902786-004 back 1.0ml		18 Aug 109 12::
50	49	08170950.d	1.	CCV 1500ng/ml S21-08170901		18 Aug 109 12::
51	50	08170951.d	1.	P0902786-005 back 1.0ml		18 Aug 109 12::
52	50	08170952.d	1.	P0902786-005dup back 1.0ml		18 Aug 109 12::
53	51	08170953.d	1.	P0902786-006 back 1.0ml		18 Aug 109 12::
54	52	08170954.d	1.	P0902786-007 back 1.0ml		18 Aug 109 12::
55	53	08170955.d	1.	P0902786-008 back 1.0ml		18 Aug 109 12::
56	54	08170956.d	1.	P0902786-009 back 1.0ml		18 Aug 109 12::
57	55	08170957.d	1.	P0902786-010 back 1.0ml		18 Aug 109 12::

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Injection Log

Directory: j:\lc01\data\to11\2009_08\17

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
58	56	08170958.d	1.	P0902786-011 back 1.0ml		18 Aug 109 12::
59	57	08170959.d	1.	P0902786-012 back 1.0ml		18 Aug 109 12::
60	58	08170960.d	1.	P0902786-013 back 1.0ml		18 Aug 109 12::
61	59	08170961.d	1.	CCV 1500ng/ml S21-08170901		18 Aug 109 12::
62	60	08170962.d	1.	ACN blk lot CY023		18 Aug 109 12::
63	61	08170963.d	1.	MB front lot 6009/6097 1.0ml		18 Aug 109 12::
64	62	08170964.d	1.	MB back lot 6009/6097 1.0ml		18 Aug 109 12::
65	63	08170965.d	1.	P0902786-014 back 1.0ml		18 Aug 109 12::
66	64	08170966.d	1.	P0902786-015 back 1.0ml		18 Aug 109 12::
67	65	08170967.d	1.	P0902786-016 back 1.0ml		18 Aug 109 12::
68	66	08170968.d	1.	P0902786-017 back 1.0ml		18 Aug 109 12::
69	67	08170969.d	1.	P0902786-018 back 1.0ml		18 Aug 109 12::
70	68	08170970.d	1.	P0902786-019 back 1.0ml		18 Aug 109 12::
71	69	08170971.d	1.	P0902786-020 back 1.0ml		18 Aug 109 12::
72	70	08170972.d	1.	P0902800-001 back 1.0ml		18 Aug 109 12::
73	71	08170973.d	1.	P0902800-002 back 1.0ml		18 Aug 109 12::
74	72	08170974.d	1.	P0902800-003 back 1.0ml		18 Aug 109 12::
75	73	08170975.d	1.	CCV 1500ng/ml S21-08170902		18 Aug 109 12::
76	74	08170976.d	1.	P0902800-004 back 1.0ml		18 Aug 109 12::
77	74	08170977.d	1.	P0902800-004dup back 1.0ml		18 Aug 109 12::
78	75	08170978.d	1.	P0902800-005 back 1.0ml		18 Aug 109 13::
79	76	08170979.d	1.	P0902800-006 back 1.0ml		18 Aug 109 13::
80	77	08170980.d	1.	P0902800-007 back 1.0ml		18 Aug 109 13::
81	78	08170981.d	1.	P0902800-008 back 1.0ml		18 Aug 109 13::
82	79	08170982.d	1.	P0902800-009 back 1.0ml		18 Aug 109 13::
83	80	08170983.d	1.	P0902800-010 back 1.0ml		18 Aug 109 13::
84	81	08170984.d	1.	P0902800-011 back 1.0ml		18 Aug 109 13::
85	82	08170985.d	1.	P0902800-012 back 1.0ml		18 Aug 109 13::
86	83	08170986.d	1.	CCV 1500ng/ml S21-08170902		18 Aug 109 13::
87	1	08170987.d	1.	ACN Blk lot CY023		18 Aug 109 13::
88	2	08170988.d	1.	MB front lot 6009/6097 1.0ml		18 Aug 109 13::
89	3	08170989.d	1.	MB back lot 6009/6097 1.0ml		18 Aug 109 13::
90	4	08170990.d	1.	P0902770-001 front 1.0ml		18 Aug 109 12::
91	5	08170991.d	1.	P0902770-002 front 1.0ml		18 Aug 109 12::
92	6	08170992.d	1.	P0902770-003 front 1.0ml		18 Aug 109 12::
93	7	08170993.d	1.	P0902770-004 front 1.0ml		18 Aug 109 12::
94	8	08170994.d	1.	P0902770-005 front 1.0ml		18 Aug 109 12::
95	9	08170995.d	1.	P0902770-006 front 1.0ml		18 Aug 109 12::
96	10	08170996.d	1.	P0902770-007 front 1.0ml		18 Aug 109 12::
97	11	08170997.d	1.	P0902770-008 front 1.0ml		18 Aug 109 12::
98	12	08170998.d	1.	P0902770-009 front 1.0ml		18 Aug 109 12::
99	13	08170999.d	1.	P0902770-010 front 1.0ml		18 Aug 109 12::
100	14	08171000.d	1.	ACN Wash		18 Aug 109 12::
101	15	08171001.d	1.	CCV 1500ng/ml S21-08170902		18 Aug 109 12::
102	16	08171002.d	1.	P0902770-011 front 1.0ml		18 Aug 109 12::
103	16	08171003.d	1.	P0902770-011dup front 1.0ml		18 Aug 109 12::
104	17	08171004.d	1.	P0902770-012 front 1.0ml		18 Aug 109 12::
105	18	08171005.d	1.	P0902770-013 front 1.0ml		18 Aug 109 12::
106	19	08171006.d	1.	P0902772-001 front 1.0ml		18 Aug 109 12::
107	20	08171007.d	1.	P0902772-002 front 1.0ml		18 Aug 109 12::
108	21	08171008.d	1.	P0902772-003 front 1.0ml		18 Aug 109 12::
109	22	08171009.d	1.	P0902772-004 front 1.0ml		18 Aug 109 12::
110	23	08171010.d	1.	P0902772-005 front 1.0ml		18 Aug 109 12::
111	24	08171011.d	1.	P0902772-006 front 1.0ml		18 Aug 109 12::
112	25	08171012.d	1.	ACN wash		18 Aug 109 12::
113	26	08171013.d	1.	CCV 1500ng/ml S21-08170902		18 Aug 109 12::
114	27	08171014.d	1.	ACN blk lot CY023		18 Aug 109 12::

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Injection Log

Directory: j:\lc01\data\to11\2009_08\17

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
115	28	08171015.d	1.	MB front lot 6009/6097 1.0ml		18 Aug 109 12::
116	29	08171016.d	1.	MB back lot 6009/6097 1.0ml		18 Aug 109 12::
117	30	08171017.d	1.	P0902772-007 front 1.0ml		18 Aug 109 12::
118	31	08171018.d	1.	P0902772-008 front 1.0ml		18 Aug 109 12::
119	32	08171019.d	1.	P0902772-009 front 1.0ml		18 Aug 109 12::
120	33	08171020.d	1.	P0902772-010 front 1.0ml		18 Aug 109 12::
121	34	08171021.d	1.	P0902772-011 front 1.0ml		18 Aug 109 12::
122	35	08171022.d	1.	P0902772-012 front 1.0ml		18 Aug 109 12::
123	36	08171023.d	1.	P0902786-001 front 1.0ml		18 Aug 109 12::
124	37	08171024.d	1.	P0902786-002 front 1.0ml		18 Aug 109 12::
125	38	08171025.d	1.	P0902786-003 front 1.0ml		18 Aug 109 12::
126	39	08171026.d	1.	P0902786-004 front 1.0ml		18 Aug 109 13::
127	40	08171027.d	1.	ACN Wash		18 Aug 109 13::
128	41	08171028.d	1.	CCV 1500ng/ml S21-08170902		18 Aug 109 13::
129	42	08171029.d	1.	P0902786-005 front 1.0ml		18 Aug 109 13::
130	42	08171030.d	1.	P0902786-005dup front 1.0ml		18 Aug 109 13::
131	43	08171031.d	1.	P0902786-006 front 1.0ml		18 Aug 109 13::
132	44	08171032.d	1.	P0902786-007 front 1.0ml		18 Aug 109 13::
133	45	08171033.d	1.	P0902786-008 front 1.0ml		18 Aug 109 13::
134	46	08171034.d	1.	P0902786-009 front 1.0ml		19 Aug 109 13::
135	47	08171035.d	1.	P0902786-010 front 1.0ml		19 Aug 109 13::
136	48	08171036.d	1.	P0902786-011 front 1.0ml		19 Aug 109 13::
137	49	08171037.d	1.	P0902786-012 front 1.0ml		19 Aug 109 13::
138	50	08171038.d	1.	P0902786-013 front 1.0ml		19 Aug 109 12::
139	51	08171039.d	1.	ACN wash		19 Aug 109 12::
140	52	08171040.d	1.	CCV 1500ng/ml S21-08170902		19 Aug 109 12::
141	53	08171041.d	1.	ACN blk lot CY023		19 Aug 109 12::
142	54	08171042.d	1.	MB front lot 6009/6097 1.0ml		19 Aug 109 12::
143	55	08171043.d	1.	MB back lot 6009/6097 1.0ml		19 Aug 109 12::
144	56	08171044.d	1.	P0902786-014 front 1.0ml		19 Aug 109 12::
145	57	08171045.d	1.	P0902786-015 front 1.0ml		19 Aug 109 12::
146	58	08171046.d	1.	P0902786-016 front 1.0ml		19 Aug 109 12::
147	59	08171047.d	1.	P0902786-017 front 1.0ml		19 Aug 109 12::
148	60	08171048.d	1.	P0902786-018 front 1.0ml		19 Aug 109 12::
149	61	08171049.d	1.	P0902786-019 front 1.0ml		19 Aug 109 12::
150	62	08171050.d	1.	P0902786-020 front 1.0ml		19 Aug 109 12::
151	63	08171051.d	1.	P0902800-001 front 1.0ml		19 Aug 109 12::
152	64	08171052.d	1.	P0902800-002 front 1.0ml		19 Aug 109 12::
153	65	08171053.d	1.	P0902800-003 front 1.0ml		19 Aug 109 12::
154	66	08171054.d	1.	ACN wash		19 Aug 109 12::
155	67	08171055.d	1.	CCV 1500ng/ml S21-08180901		19 Aug 109 12::
156	68	08171056.d	1.	P0902800-004 front 1.0ml		19 Aug 109 12::
157	68	08171057.d	1.	P0902800-004dup front 1.0ml		19 Aug 109 12::
158	69	08171058.d	1.	P0902800-005 front 1.0ml		19 Aug 109 12::
159	70	08171059.d	1.	P0902800-006 front 1.0ml		19 Aug 109 12::
160	71	08171060.d	1.	P0902800-007 front 1.0ml		19 Aug 109 12::
161	72	08171061.d	1.	P0902800-008 front 1.0ml		19 Aug 109 12::
162	73	08171062.d	1.	P0902800-009 front 1.0ml		19 Aug 109 12::
163	74	08171063.d	1.	P0902800-010 front 1.0ml		19 Aug 109 12::
164	75	08171064.d	1.	P0902800-011 front 1.0ml		19 Aug 109 12::
165	76	08171065.d	1.	ACN wash		19 Aug 109 12::
166	77	08171066.d	1.	CCV 1500ng/ml S21-08180901		19 Aug 109 12::
167	78	08171067.d	1.	ACN blk lot CYo23		19 Aug 109 12::
168	79	08171068.d	1.	MB front lot 6009/6097 1.0ml		19 Aug 109 12::
169	80	08171069.d	1.	MB back lot 6009/6097 1.0ml		19 Aug 109 12::
170	81	08171070.d	1.	P0902800-013 back 1.0ml		19 Aug 109 12::
171	82	08171071.d	1.	P0902800-012 front 1.0ml		19 Aug 109 12::

Injection Log

Directory: j:\lc01\data\to11\2009_08\17

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
172	82	08171072.d	1.	P0902800-012dup front 1.0ml		19 Aug 109 12::
173	83	08171073.d	1.	P0902800-013 front 1.0ml		19 Aug 109 12::
174	84	08171074.d	1.	ACN wash		19 Aug 109 13::
175	85	08171075.d	1.	CCV 1500ng/ml S21-08180901		19 Aug 109 13::
176	1	08171076.d	1.	ACN lot CY023		19 Aug 109 13::
177	2	08171077.d	1.	P0902770-001 front 10x		19 Aug 109 13::
178	3	08171078.d	1.	P0902770-002 front 10x		19 Aug 109 13::
179	4	08171079.d	1.	P0902770-004 front 10x		19 Aug 109 13::
180	5	08171080.d	1.	P0902770-005 front 10x		19 Aug 109 13::
181	6	08171081.d	1.	P0902772-007 front 10x		19 Aug 109 13::
182	7	08171082.d	1.	P0902772-008 front 10x		19 Aug 109 13::
183	8	08171083.d	1.	P0902772-011 front 10x		19 Aug 109 13::
184	9	08171084.d	1.	P0902772-012 front 10x		19 Aug 109 13::
185	10	08171085.d	1.	P0902771-007 front 10x		19 Aug 109 13::
186	11	08171086.d	1.	CCV 1500ng/ml S21-08180901		19 Aug 109 12::