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

LABORATORY REPORT

September 21, 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 25, 2009. For your reference, these analyses have been assigned our service request number P0902942.

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 335 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: P0902942
Project: 16512

CASE NARRATIVE

The samples were received intact under chain of custody on August 25, 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.

Aldehyde Analysis

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

The minimum control criterion for 2,5-Dimethylhexaldehyde and Crotonaldehyde were outside the continuing calibration verification (CCV) method requirements. The analytes were not detected in the samples; therefore, the method reporting limits (MRL) have been raised to account for the bias. Additionally, the data is flagged with the appropriate data qualifier.

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, Inc. **Service Request:** P0902942
Project: 16512

SAMPLE CROSS-REFERENCE

SAMPLE #	CLIENT SAMPLE ID	DATE	TIME
P0902942-001	101469	8/24/09	00:00
P0902942-002	101470	8/24/09	00:00
P0902942-003	101471	8/24/09	00:00
P0902942-004	101472	8/24/09	00:00
P0902942-005	101473	8/24/09	00:00
P0902942-006	101474	8/24/09	00:00

Environmental Health & Engineering, Inc.

CHAIN OF CUSTODY FORM

DATE: 8/24/09

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

P0902942

TO: Columbia Analytical

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	101469	Sorbent - Air	EPA TO-11 Full List	104.04 L
2	101470	"	"	104.04 L
3	101471	"	"	100.5 L
4	101472	"	"	101.5 L
5	101473	"	"	104.54 L
6	101474	"	"	Ø L FB

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: Tuan Truong of Environmental Health & Engineering, Inc. Date: 8/24/09

Received by: _____ of (company name) CAS Date: 8/25/09 0945

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

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: _____

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, Inc. Work order: P0902942

Project: 16512

Sample(s) received on: 08/25/09 Date opened: 08/25/09 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 3 °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
P0902942-001.01	Silica Gel DNPH Tube					
P0902942-002.01	Silica Gel DNPH Tube					
P0902942-003.01	Silica Gel DNPH Tube					
P0902942-004.01	Silica Gel DNPH Tube					
P0902942-005.01	Silica Gel DNPH Tube					
P0902942-006.01	Silica Gel DNPH Tube					

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

Chain of Custody is missing time collected

*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: **101469**
Client Project ID: **16512**

CAS Project ID: P0902942
CAS Sample ID: P0902942-001

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: 8/24/09
Date Received: 8/25/09
Date Analyzed: 8/26-27/09
Desorption Volume: 1.0 ml
Volume Sampled: 104.04 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**2,300**	**22**	0.96	**18**	0.78	
75-07-0	**Acetaldehyde**	**4,600**	**44**	0.96	**25**	0.53	
123-38-6	**Propionaldehyde**	**590**	**5.7**	0.96	**2.4**	0.40	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.96	ND	0.34	
123-72-8	**Butyraldehyde**	**1,000**	**10**	0.96	**3.4**	0.33	
100-52-7	**Benzaldehyde**	**960**	**9.2**	0.96	**2.1**	0.22	
590-86-3	**Isovaleraldehyde**	**150**	**1.5**	0.96	**0.42**	0.27	
110-62-3	**Valeraldehyde**	**1,600**	**16**	0.96	**4.4**	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**	**4,500**	**43**	0.96	**11**	0.23	
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.

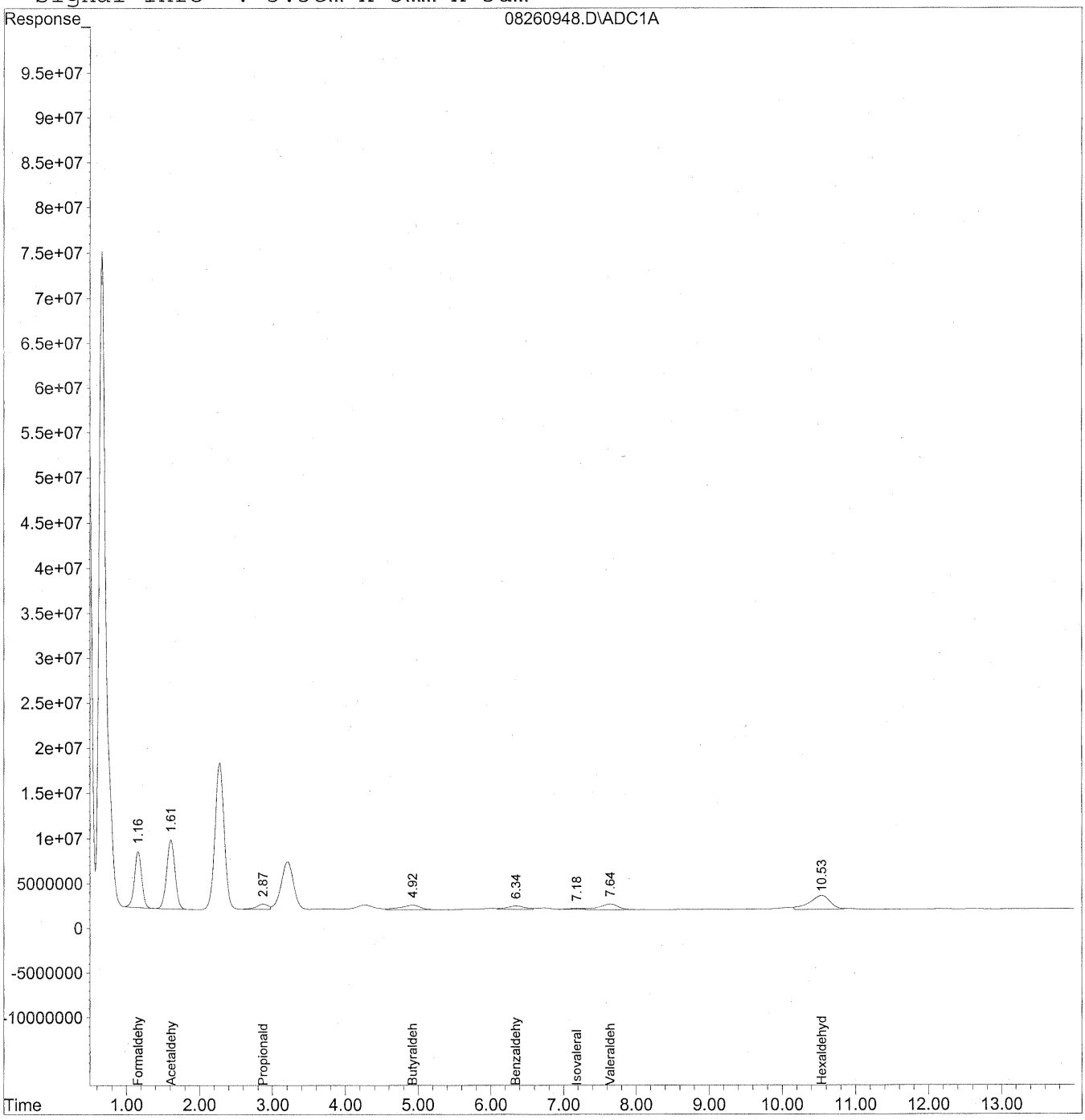
Verified By: Rc Date: 9/11/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260948.D Vial: 47
Acq On : 27 Aug 2009 4:52 am Operator: HC
Sample : P0902942-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:35 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 : Thu Aug 27 07:35:56 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



8

Quantitation Report (QT Reviewed)

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Acq On : 27 Aug 2009 4:52 am Operator: HC
Sample : P0902942-001 front 1.0ml Inst : LC 01
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IntFile : autoint1.e
Quant Time: Aug 29 17:35 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 : Thu Aug 27 07:35:56 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	425733090	2319.043	ng/ml
2)	Acetaldehyde	1.61	613879290	4377.863	ng/ml
3)	Propionaldehyde	2.87	63425776	594.458	ng/mlm
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	4.92	92427864	1046.320	ng/mlm
6)	Benzaldehyde	6.34	63013773	956.648	ng/mlm
7)	Isovaleraldehyde	7.18	12121187	154.901	ng/mlm
8)	Valeraldehyde	7.64	118710062	1614.993	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	10.53	304651143	4523.821	ng/mlm
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/mld

Quantitation Report

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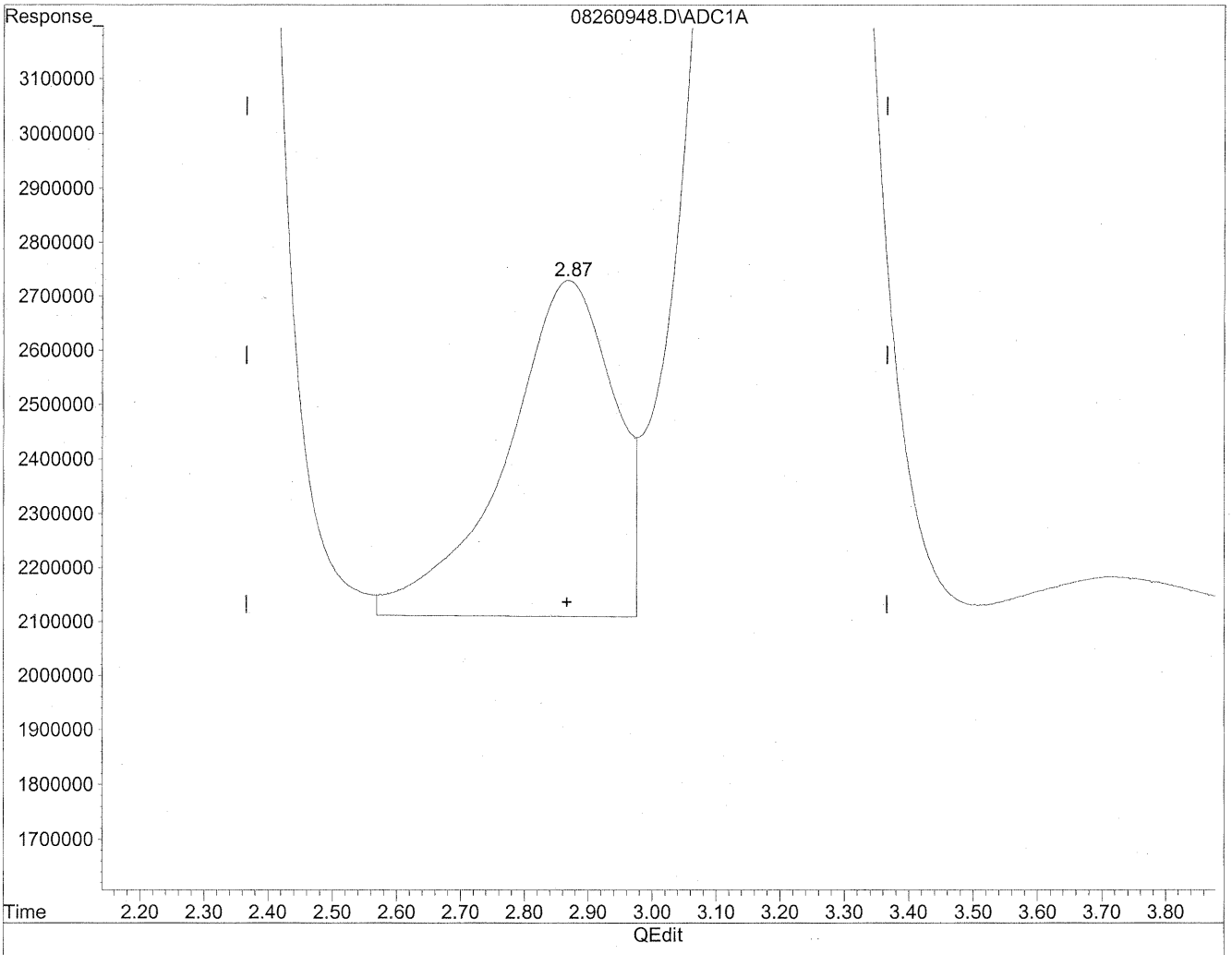
Sample : P0902942-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:36 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.87min 684.147ng/ml

response 72995223

(+) = Expected Retention Time

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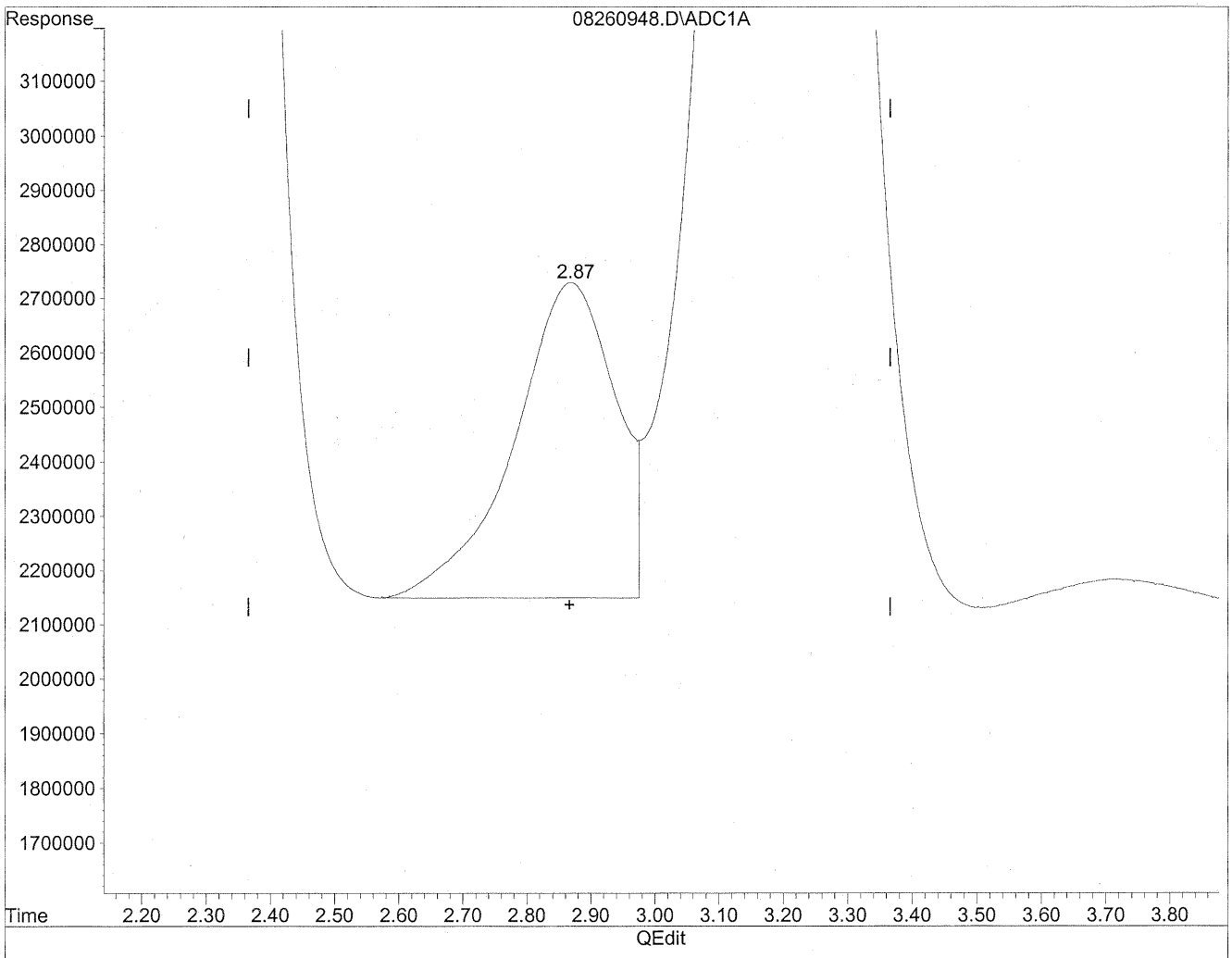

(3) Propionaldehyde

2.87min 594.458ng/ml m

response 63425776

HC 8/28/09
BC

LM 8/31/09

(+) = Expected Retention Time

Quantitation Report

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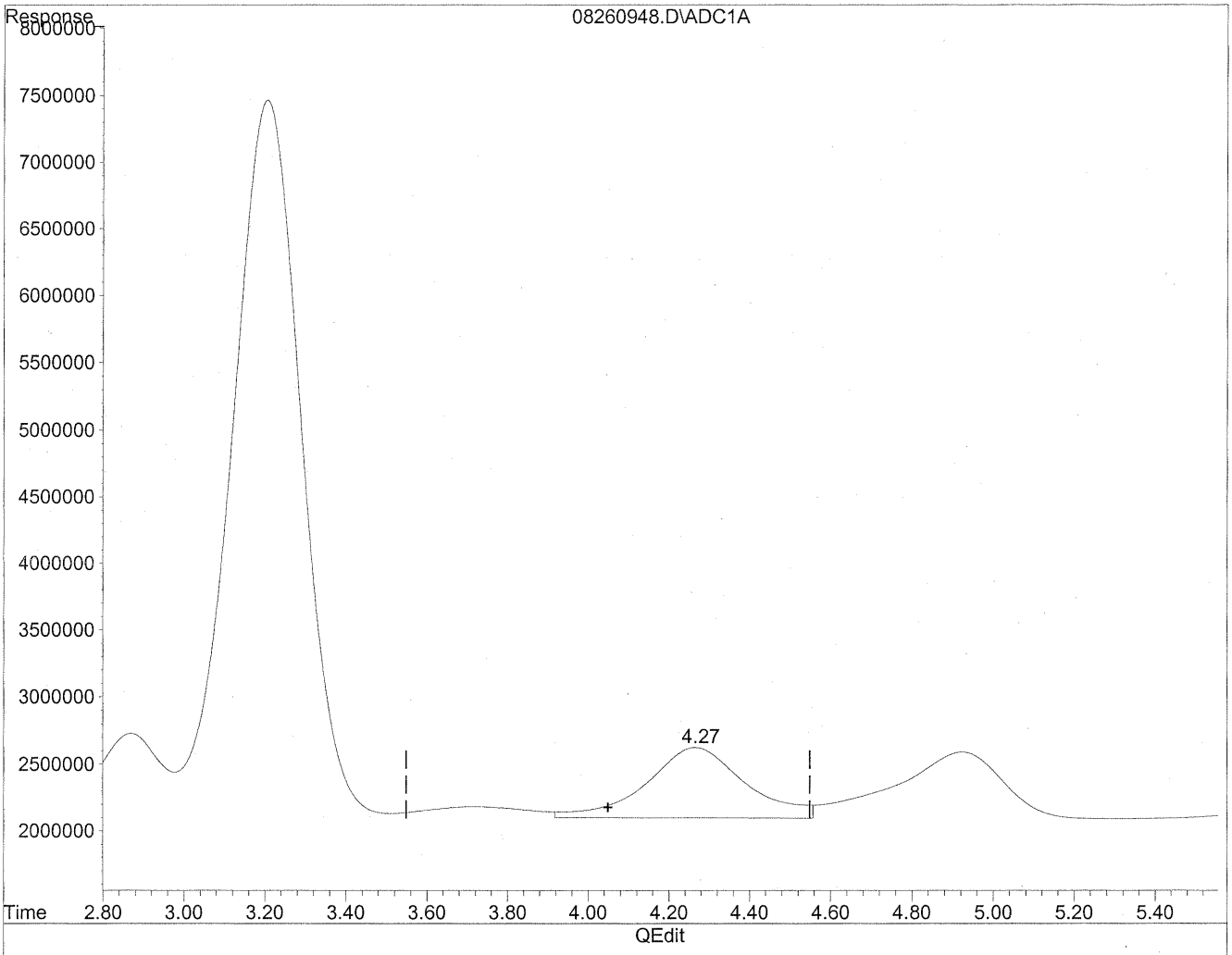

(4) Crotonaldehyde

4.26min 916.256ng/ml

response 89257265

(+) = Expected Retention Time

Quantitation Report

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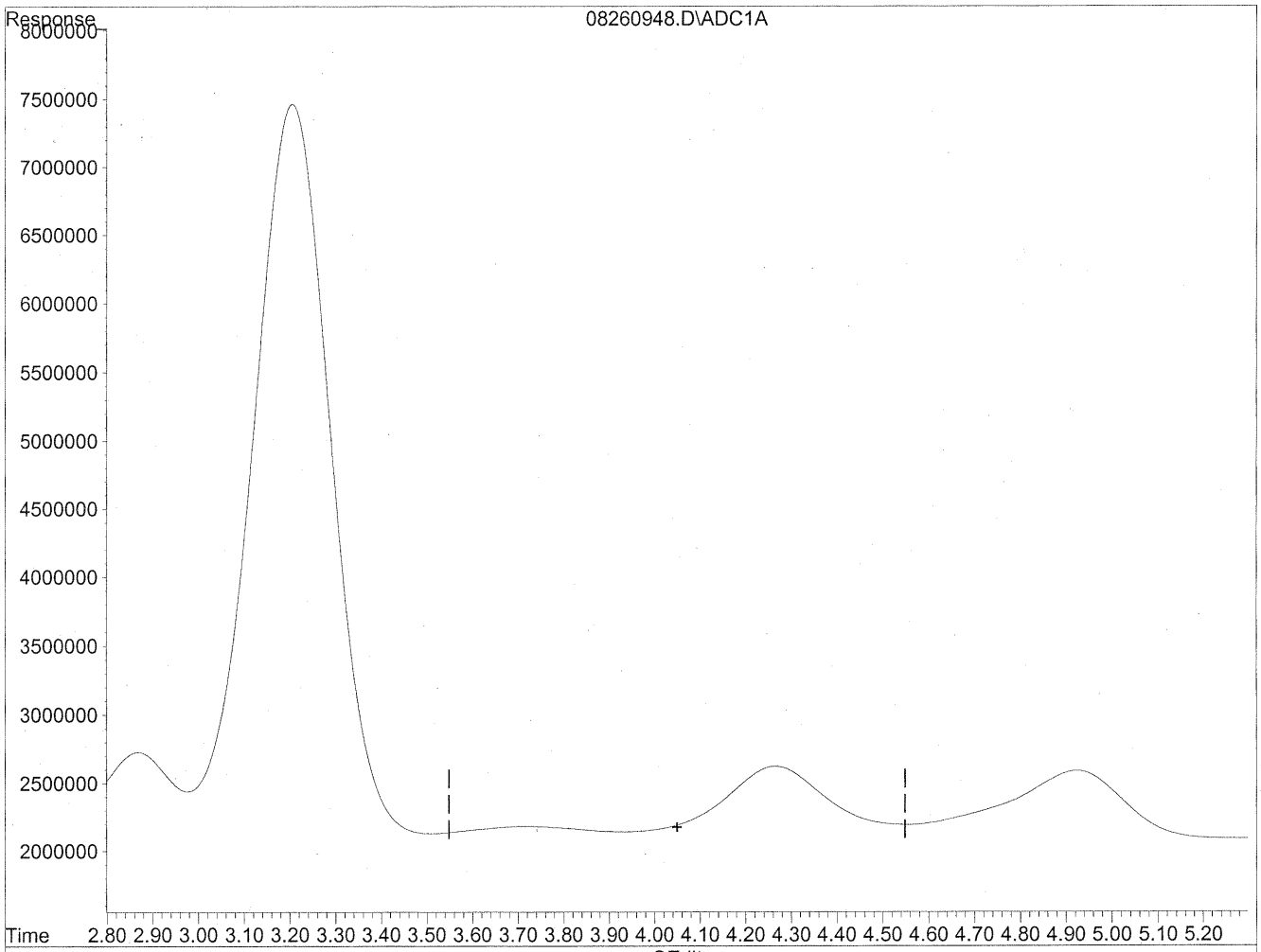

(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/31/09 WP

CW 8/31/09

(+) = Expected Retention Time

Quantitation Report

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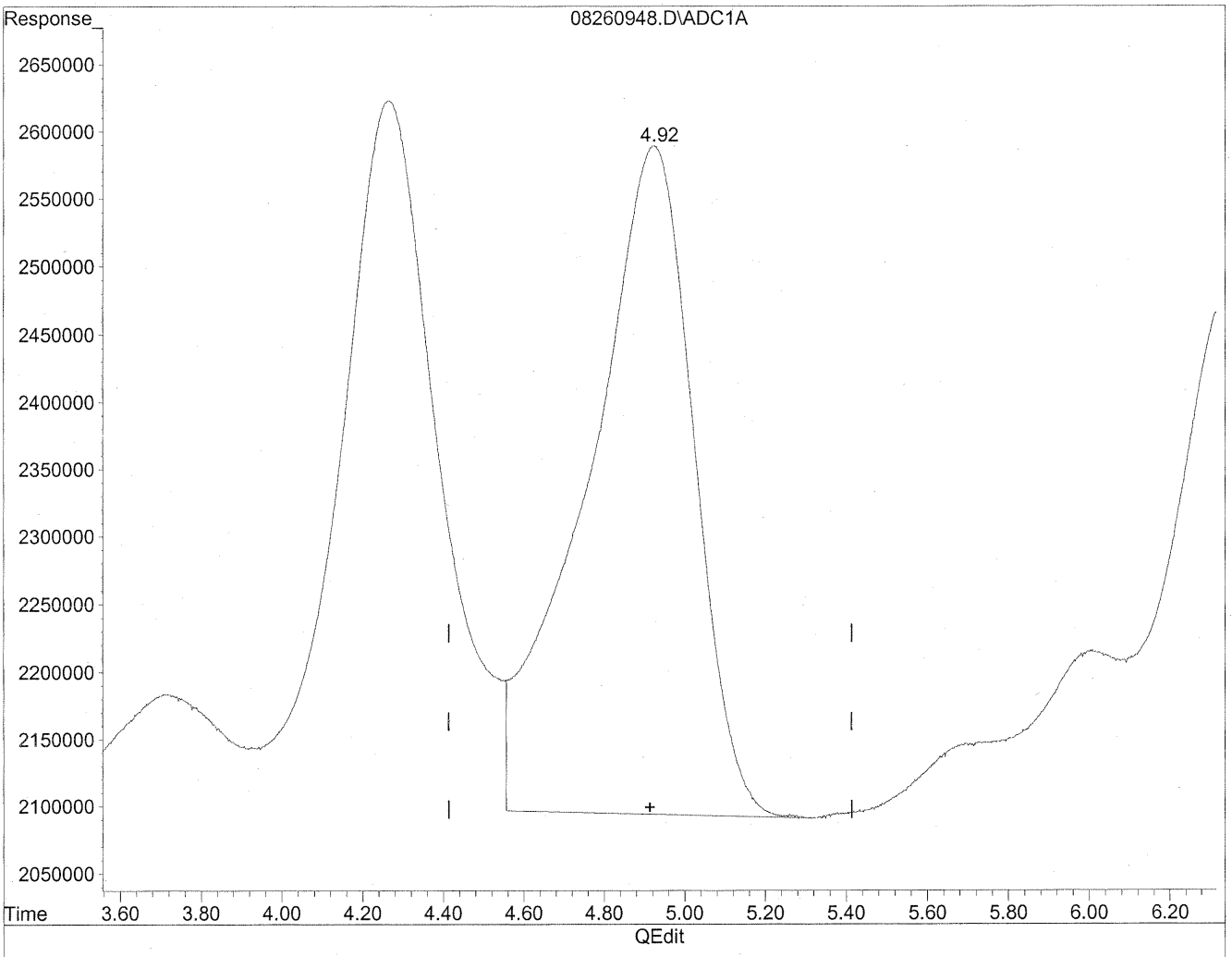

(5) Butyraldehyde

4.92min 1027.784ng/ml

response 90790468

(+) = Expected Retention Time

08260948.D TO110709.M Sat Aug 29 17:33:21 2009

Quantitation Report

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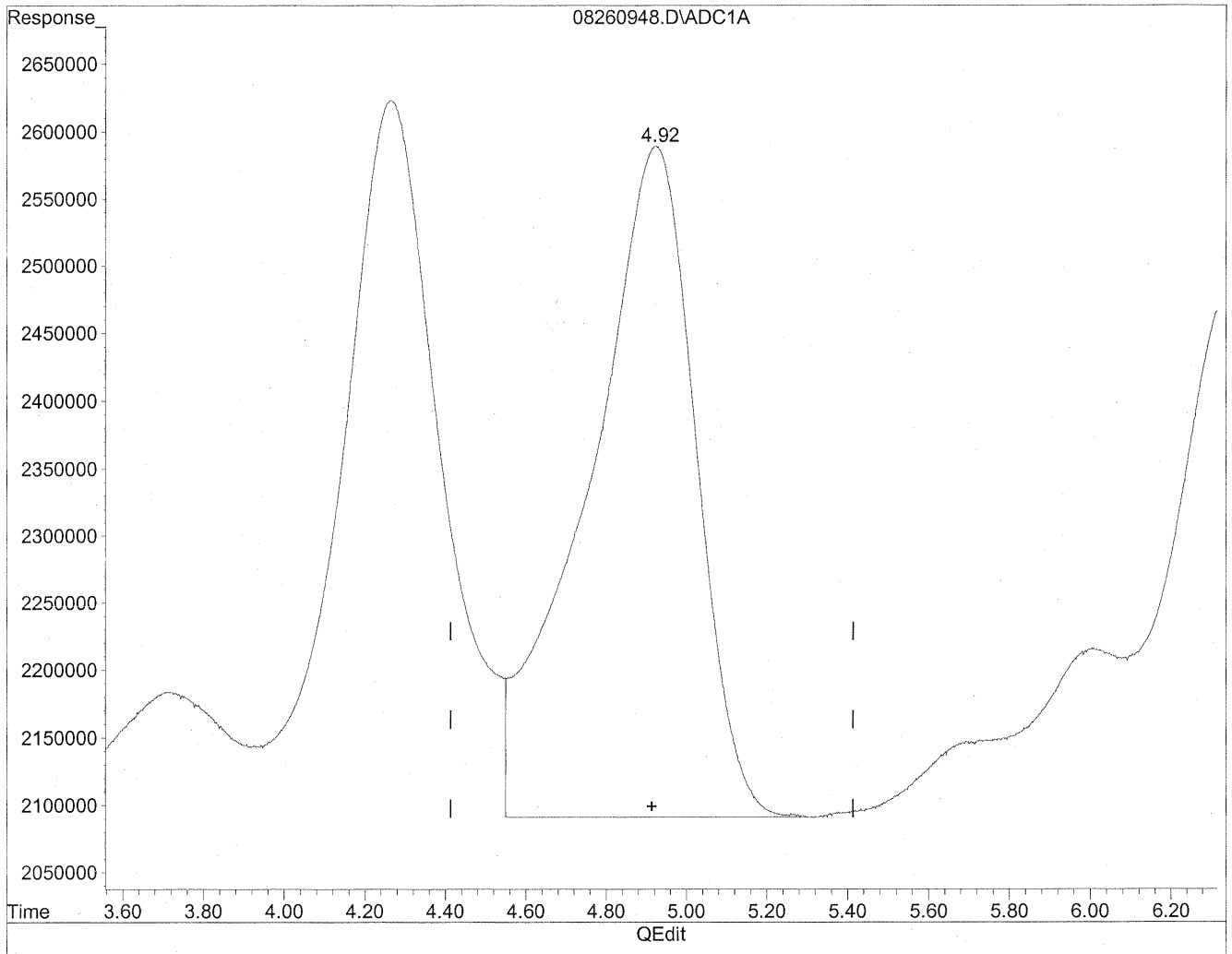

(5) Butyraldehyde

4.92min 1046.320ng/ml m

response 92427864

HC 8/31/09

RL

8/31/09

(+) = Expected Retention Time

08260948.D TO110709.M Sat Aug 29 17:33:31 2009

Quantitation Report

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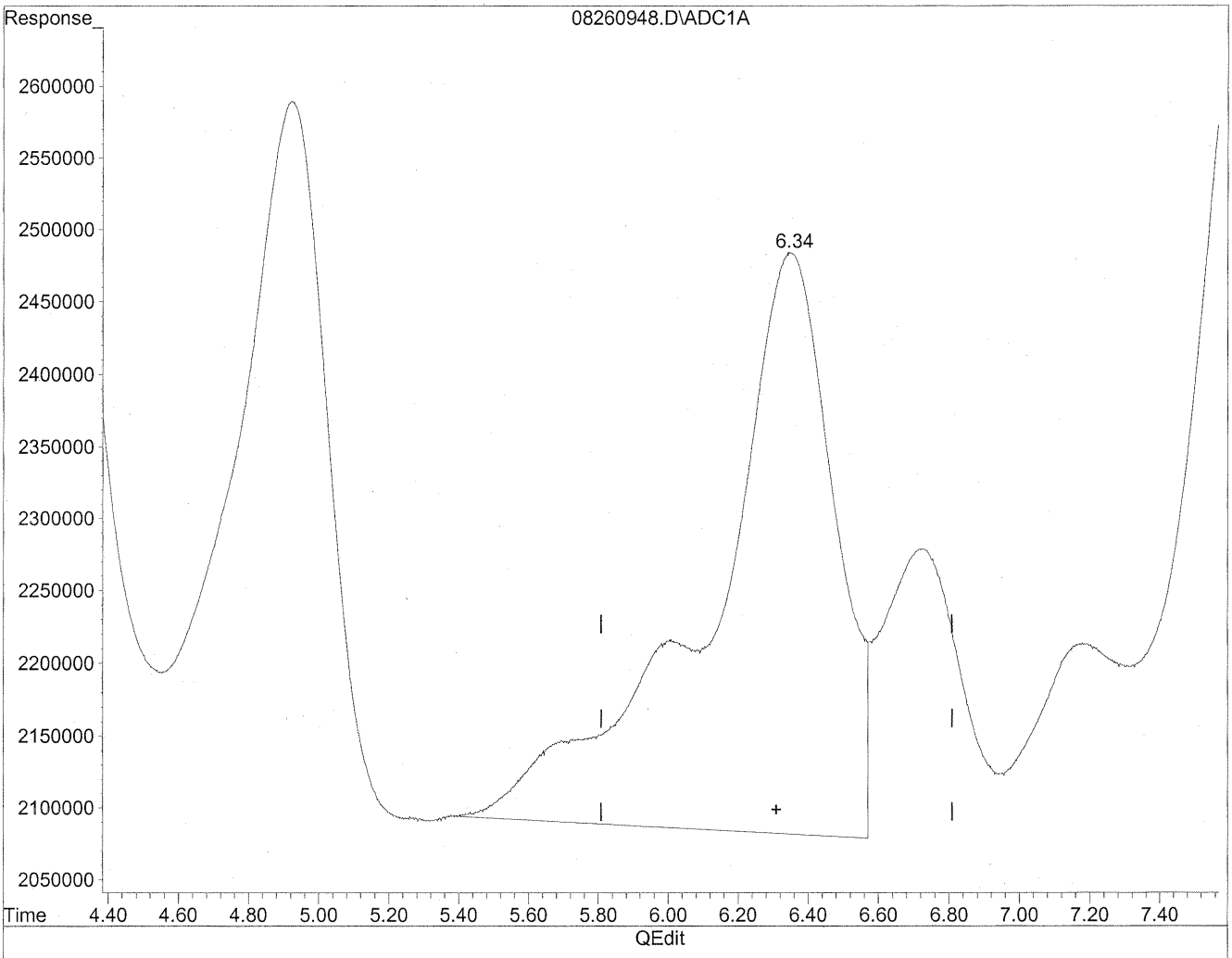

(6) Benzaldehyde

6.35min 1507.391ng/ml

response 99290821

(+) = Expected Retention Time

Quantitation Report

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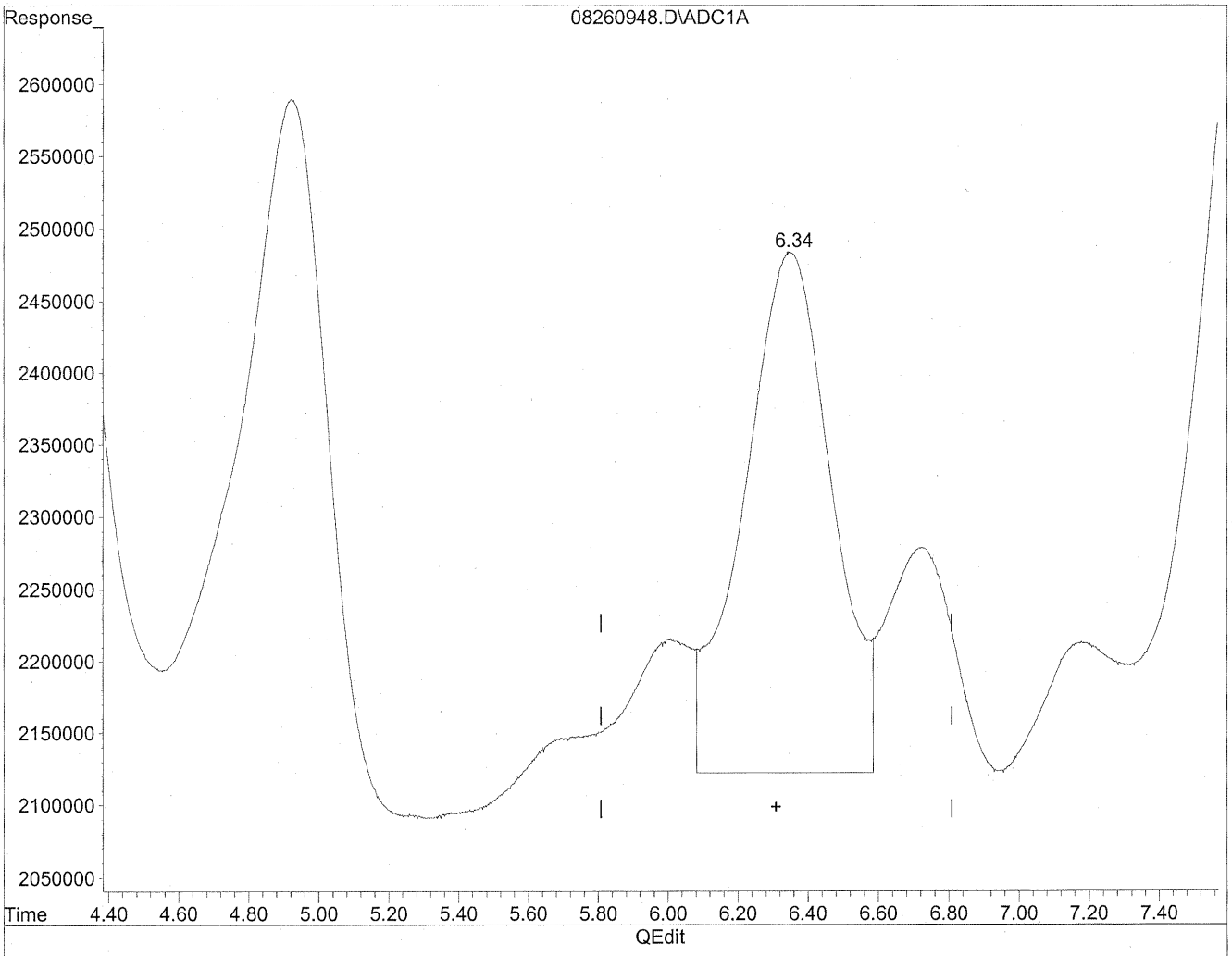

(6) Benzaldehyde

6.34min 956.648ng/ml m

response 63013773

tlc 8/31/09
BE

8/31/09

(+) = Expected Retention Time

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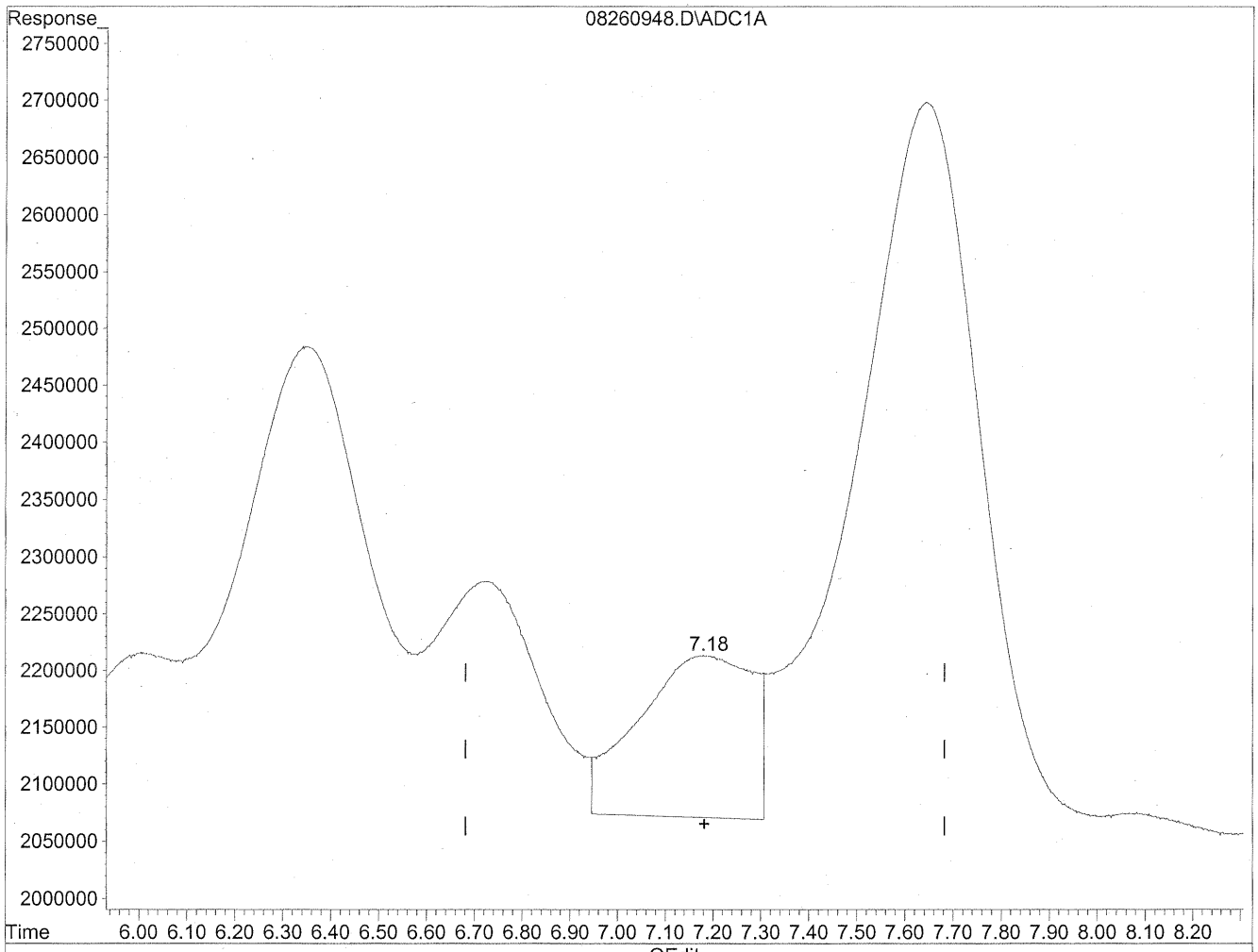

QEdit

(7) Isovaleraldehyde

7.18min 303.349ng/ml

response 23737364

(+) = Expected Retention Time

Quantitation Report

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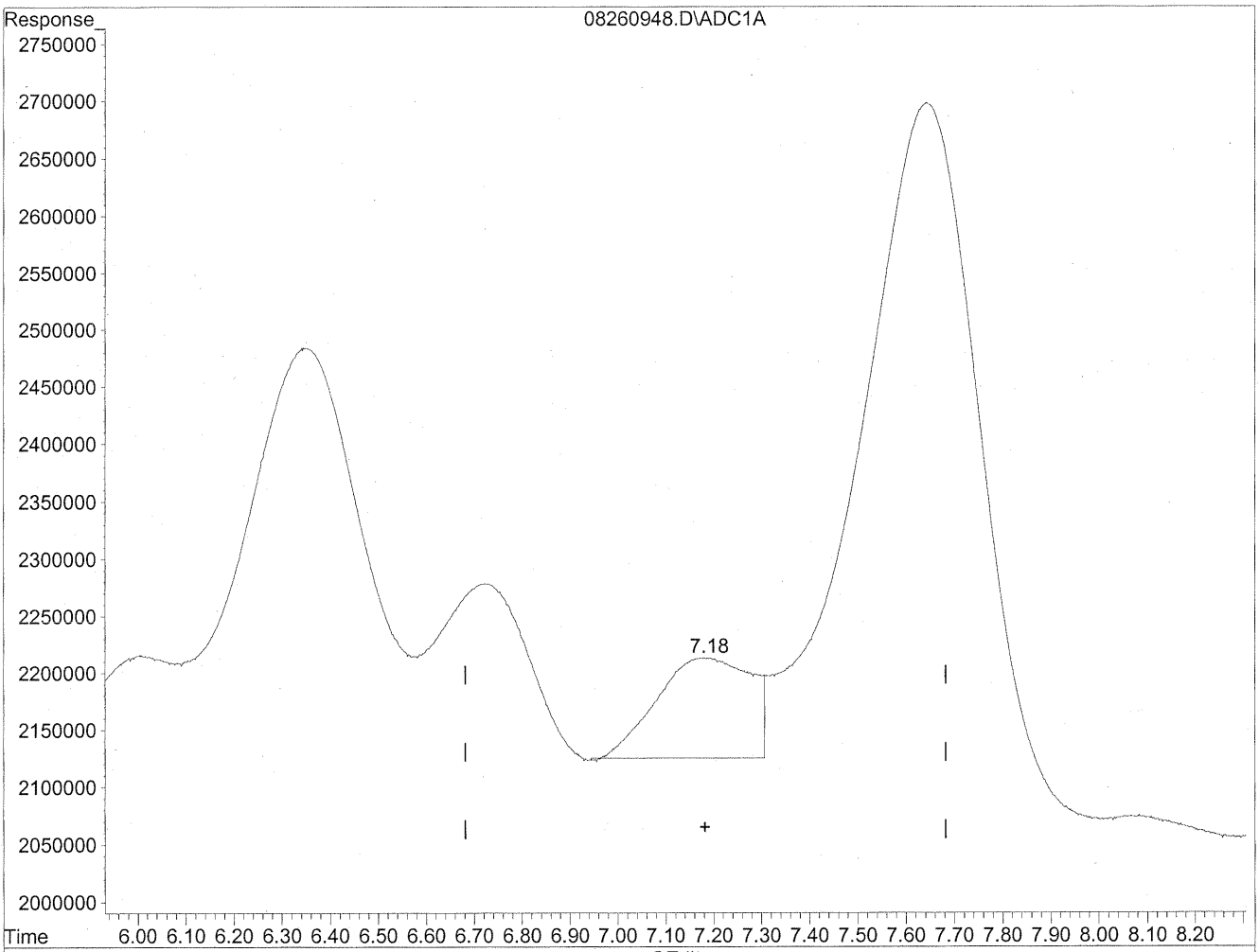

(7) Isovaleraldehyde

7.18min 154.901ng/ml m

response 12121187

(+) = Expected Retention Time

08260948.D TO110709.M Sat Aug 29 17:34:11 2009

Quantitation Report

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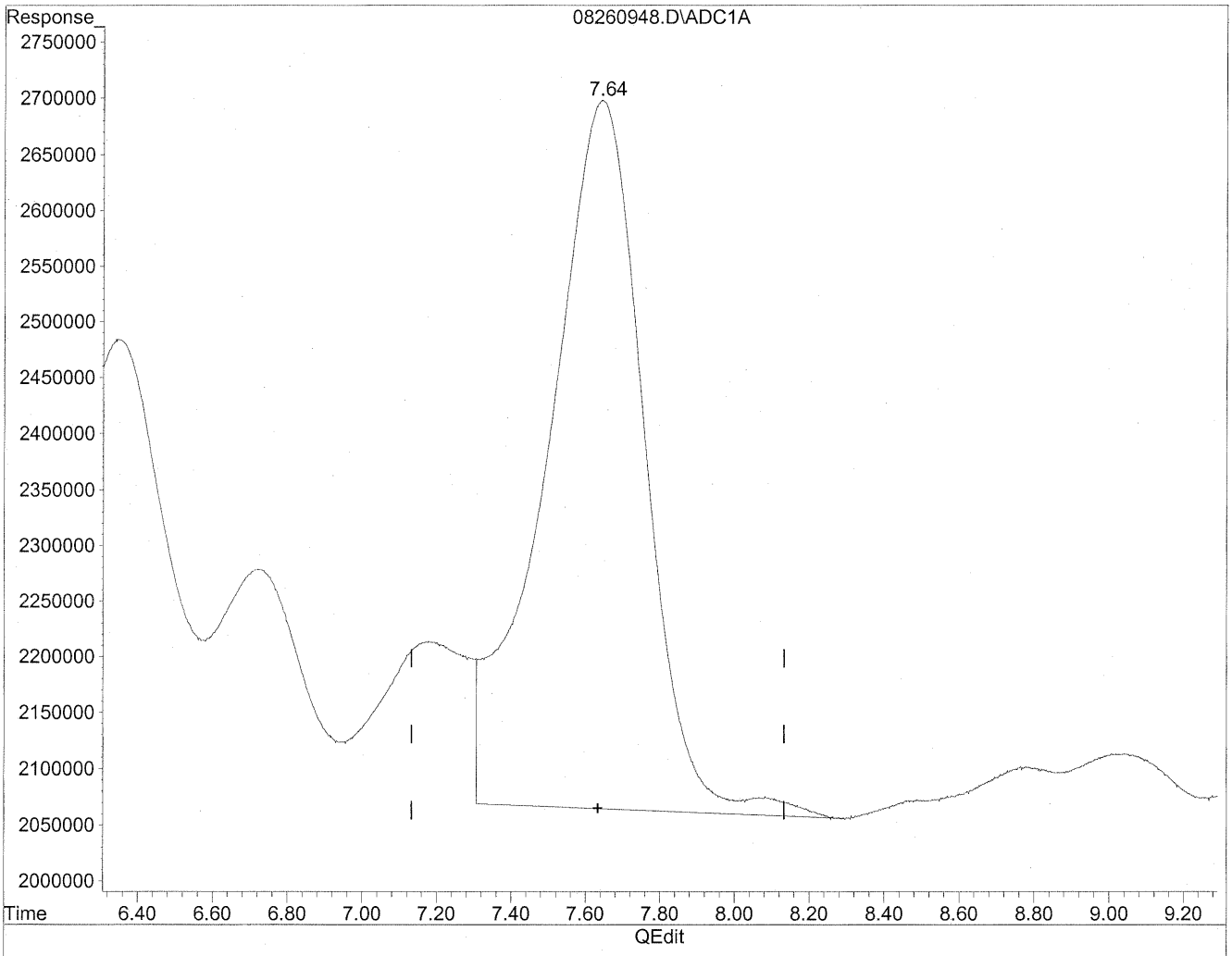
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.64min 1575.535ng/ml

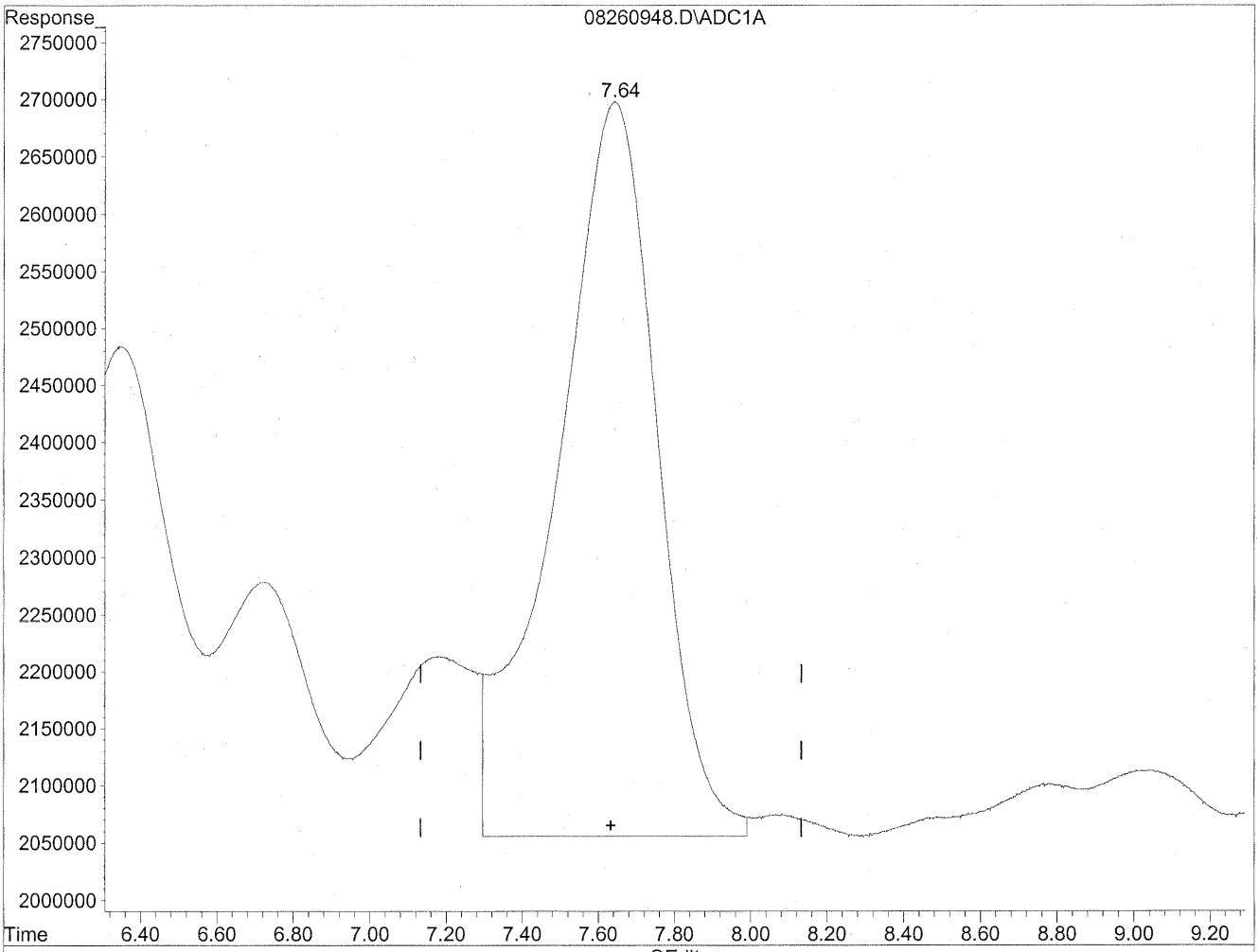
response 115809726

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260948.D Vial: 47
Acq On : 27 Aug 2009 4:52 am Operator: HC
Sample : P0902942-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:36 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.64min 1614.993ng/ml m

response 118710062

HC 8/31/09
SH/BC

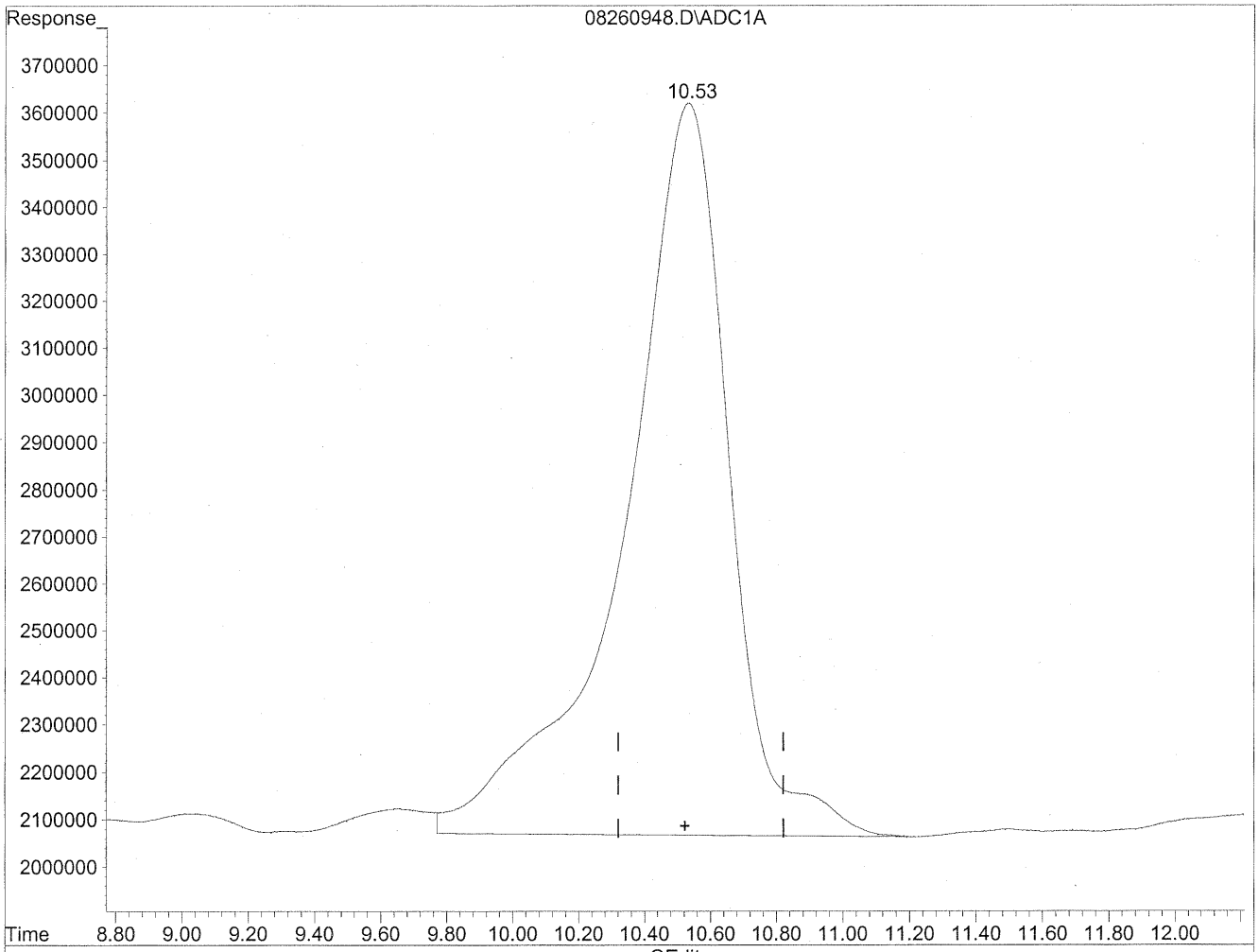
(+) = Expected Retention Time

08260948.D TO110709.M Sat Aug 29 17:34:34 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260948.D Vial: 47
Acq On : 27 Aug 2009 4:52 am Operator: HC
Sample : P0902942-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:36 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.53min 5111.334ng/ml

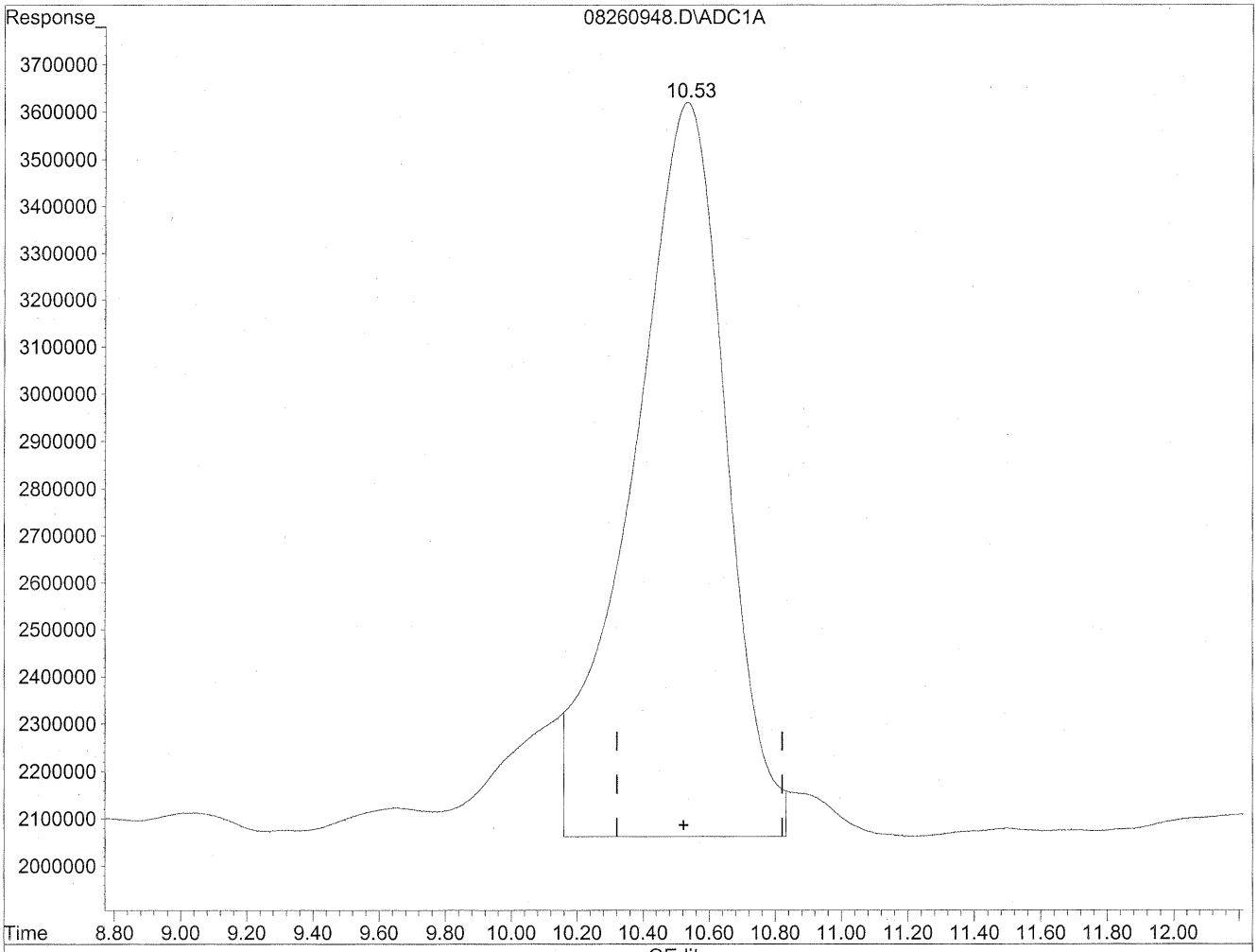
response 344216452

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260948.D Vial: 47
Acq On : 27 Aug 2009 4:52 am Operator: HC
Sample : P0902942-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:36 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.53min 4523.821ng/ml m

response 304651143

HC 8/31/09 STL, BC

8/31/09

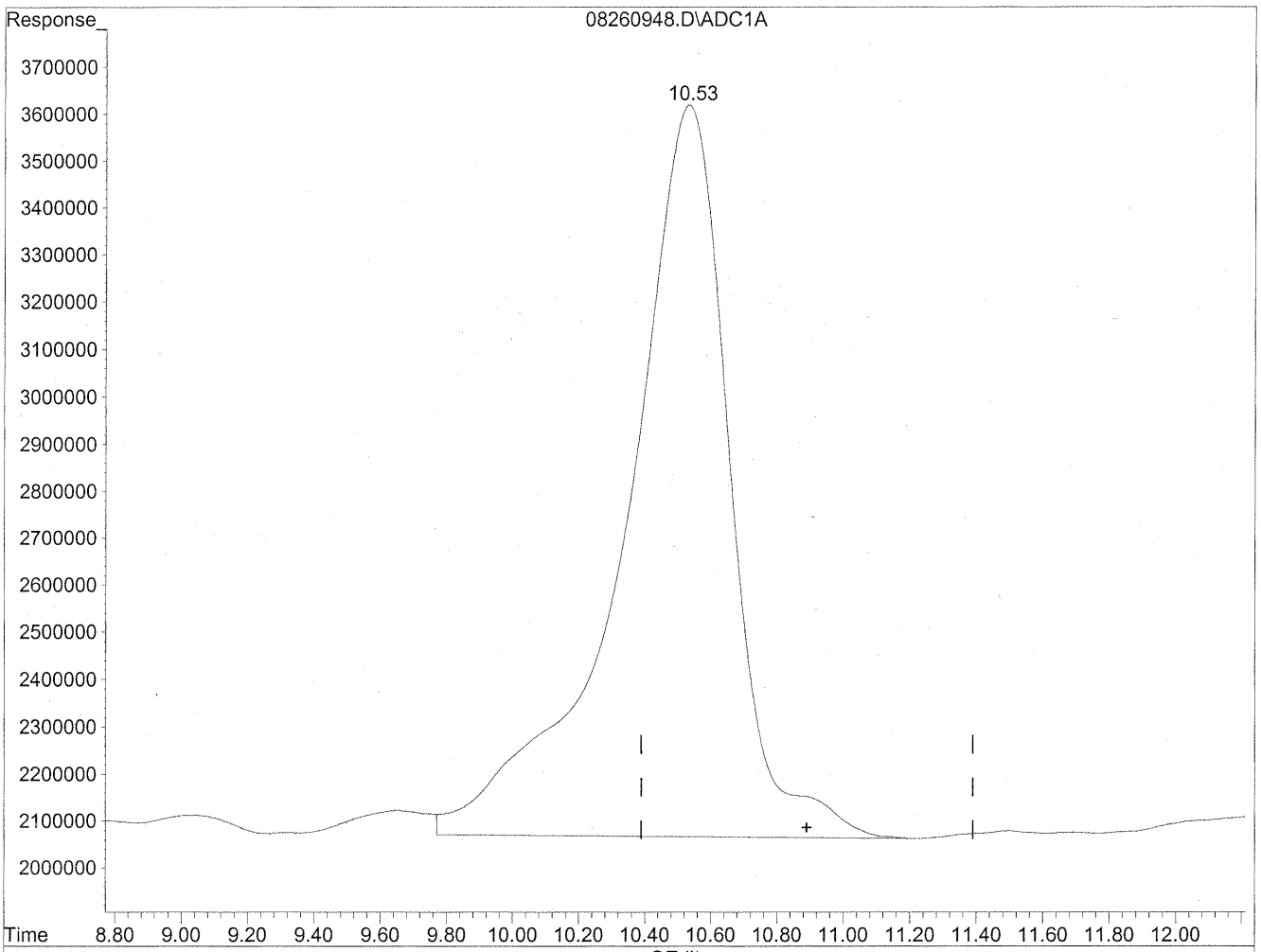
23

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260948.D Vial: 47
Acq On : 27 Aug 2009 4:52 am Operator: HC
Sample : P0902942-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:36 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.53min 7022.906ng/ml

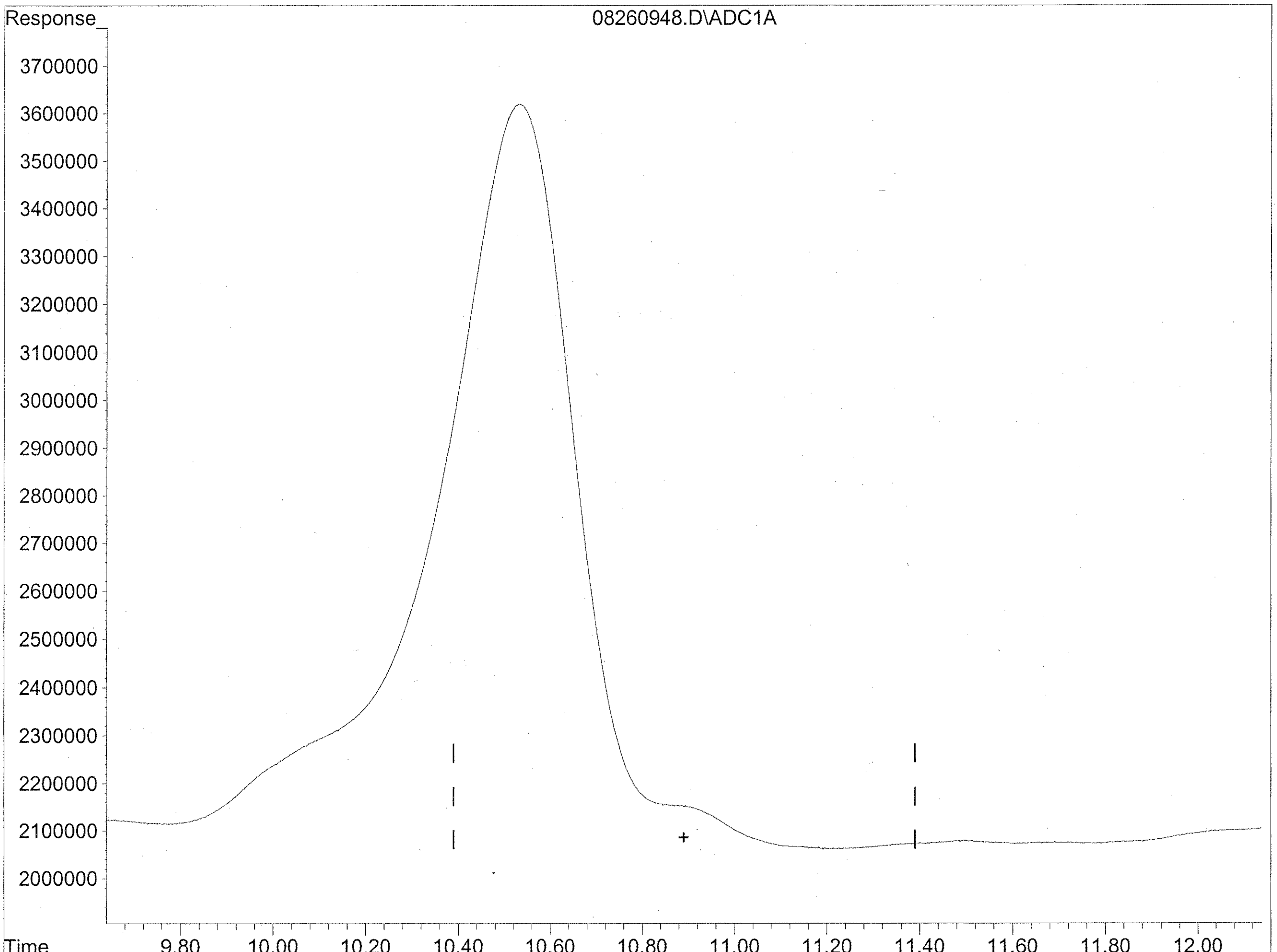
response 344216452

(+) = Expected Retention Time

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2009_08\26\08260948.D       Vial: 47
Acq On    : 27 Aug 2009   4:52 am                       Operator: HC
Sample    : P0902942-001 front 1.0ml                     Inst    : LC 01
Misc      :                                             Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Aug 27  7:36 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    : Sat Aug 29 16:33:38 2009
Response via   : Multiple Level Calibration
```



QEdit

(12) 2,5-Dimethylbenzaldehyde

0.00min   0.000ng/ml d

response   0

HC 8/31/09 ND

GW 8/31/09

25

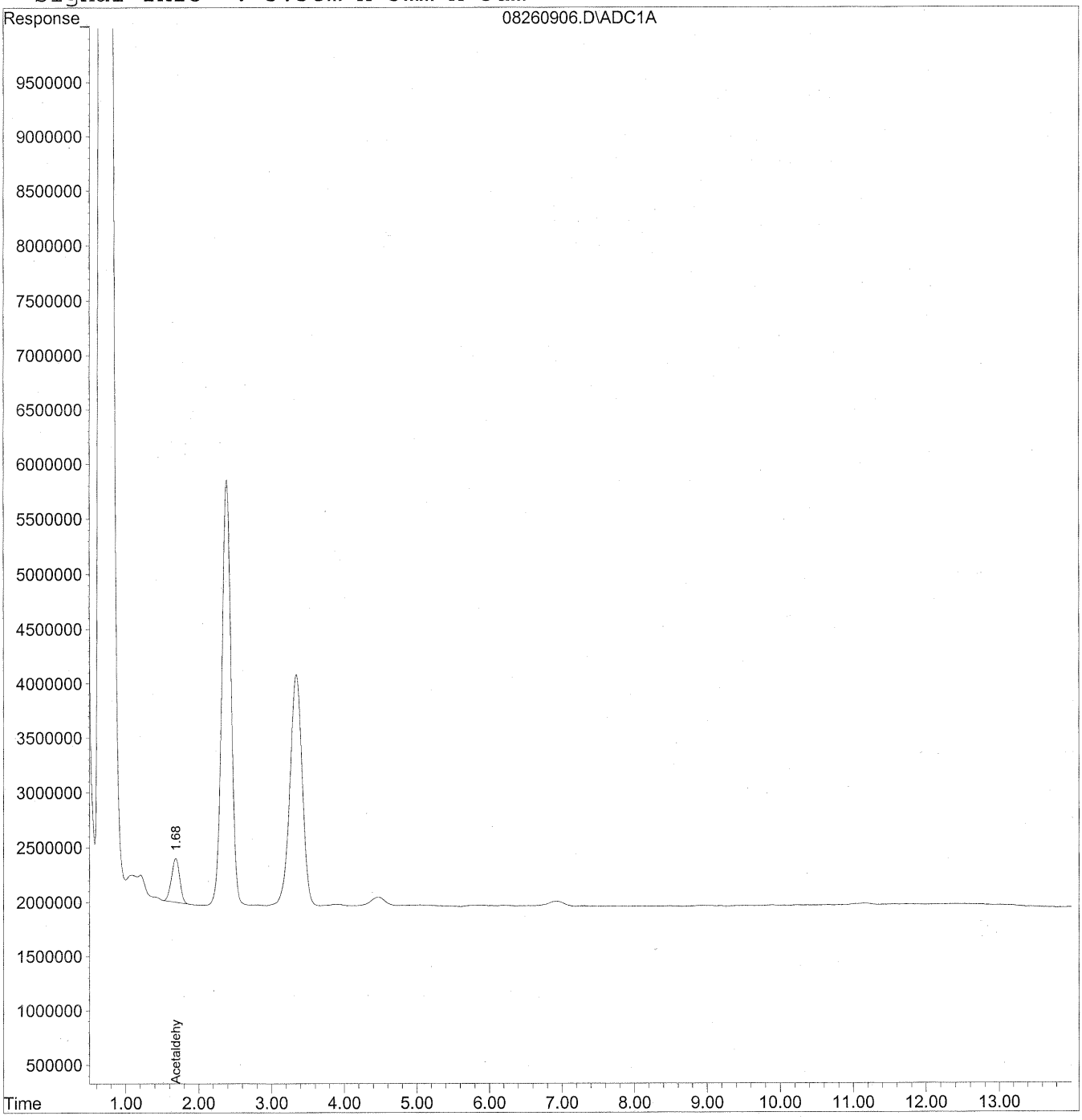
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260906.D Vial: 6
Acq On : 26 Aug 2009 6:20 pm Operator: HC
Sample : P0902942-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:46 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 : Sat Aug 29 16:33:38 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



26

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\26\08260906.D Vial: 6
Acq On : 26 Aug 2009 6:20 pm Operator: HC
Sample : P0902942-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:46 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 : Sat Aug 29 16:33:38 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

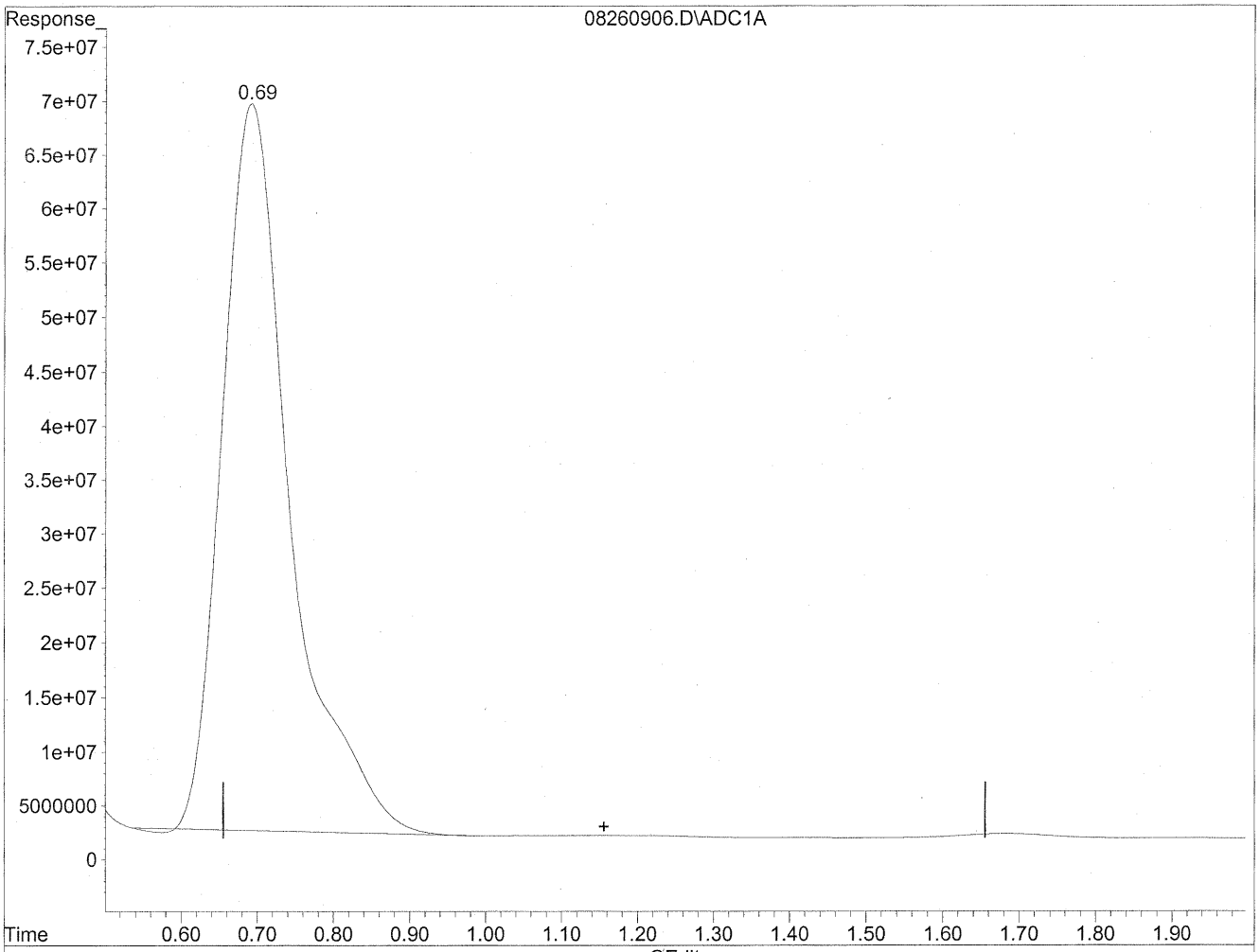
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	0.00	0	N.D.	ng/mld
2)	Acetaldehyde	1.68	32327639	230.544	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/mld
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\26\08260906.D Vial: 6
Acq On : 26 Aug 2009 6:20 pm Operator: HC
Sample : P0902942-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:45 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

0.69min 22631.270ng/ml

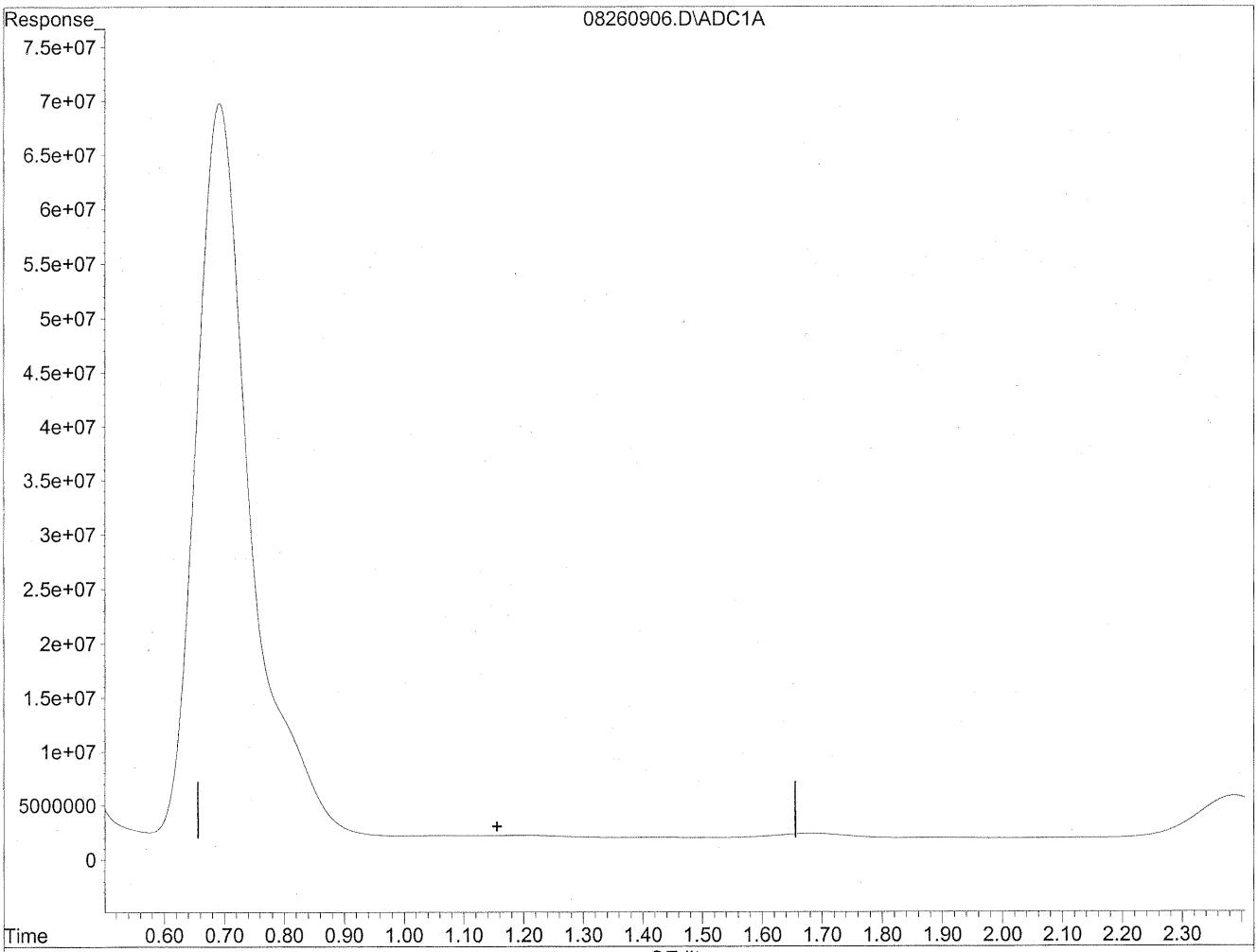
response 4154680056

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260906.D Vial: 6
Acq On : 26 Aug 2009 6:20 pm Operator: HC
Sample : P0902942-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:45 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

0.00min 0.000ng/ml d

response 0

8/30/09

8/31/09

29

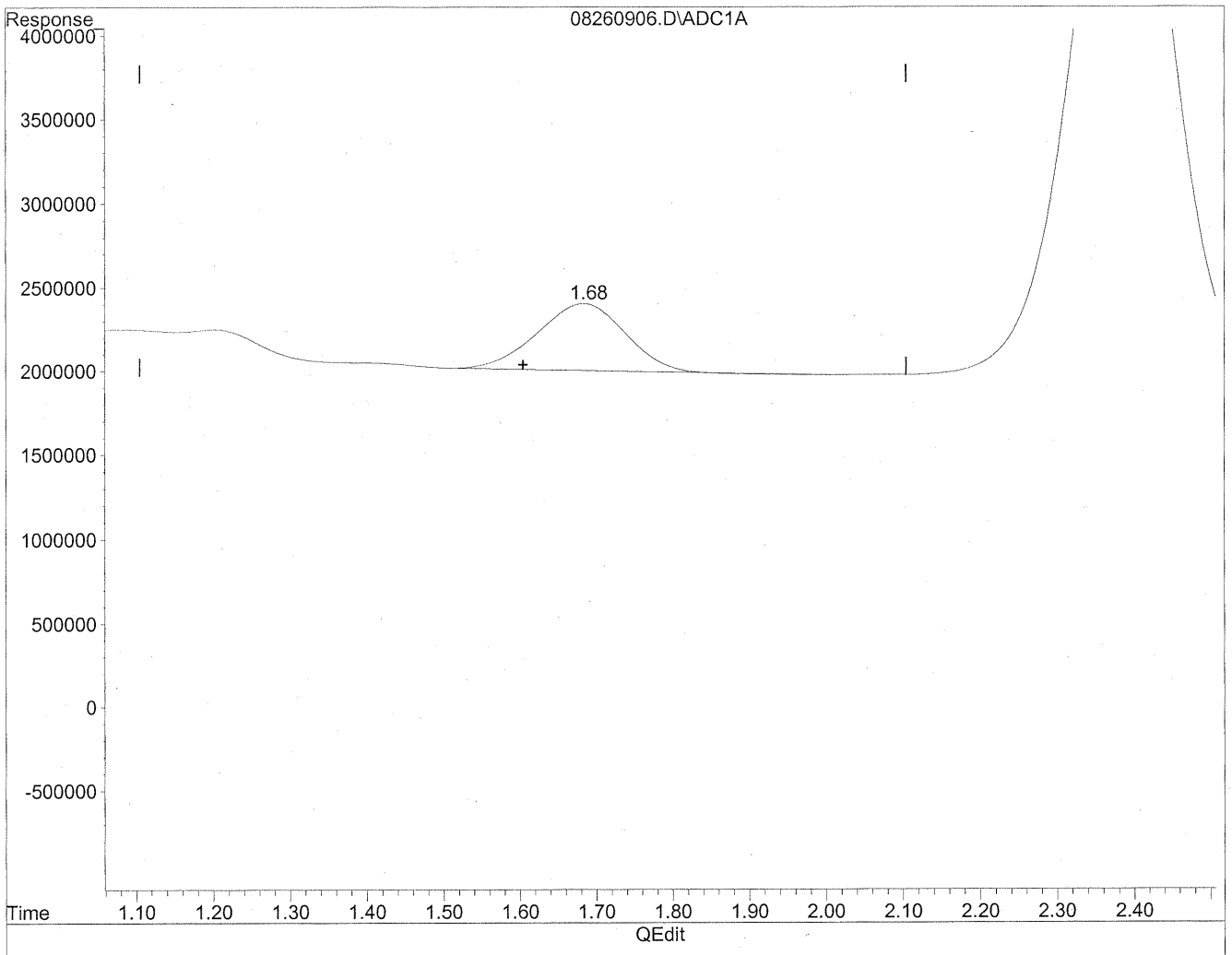
(+) = Expected Retention Time

08260906.D TO110709.M Sat Aug 29 16:45:58 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260906.D Vial: 6
Acq On : 26 Aug 2009 6:20 pm Operator: HC
Sample : P0902942-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:45 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.68min 226.192ng/ml

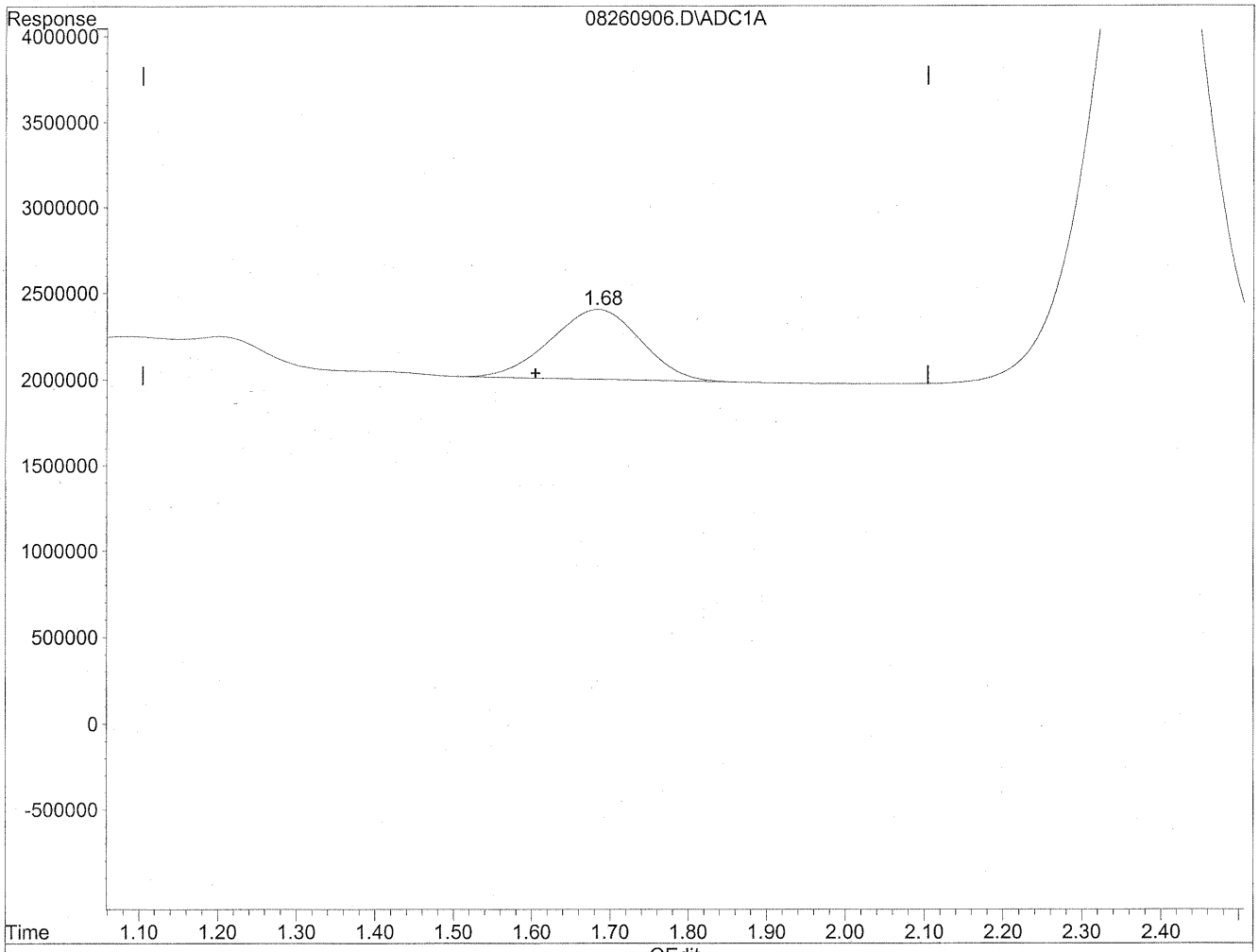
response 31717503

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260906.D Vial: 6
Acq On : 26 Aug 2009 6:20 pm Operator: HC
Sample : P0902942-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:45 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.68min 230.544ng/ml m

response 32327639

HC 8/31/09 IC

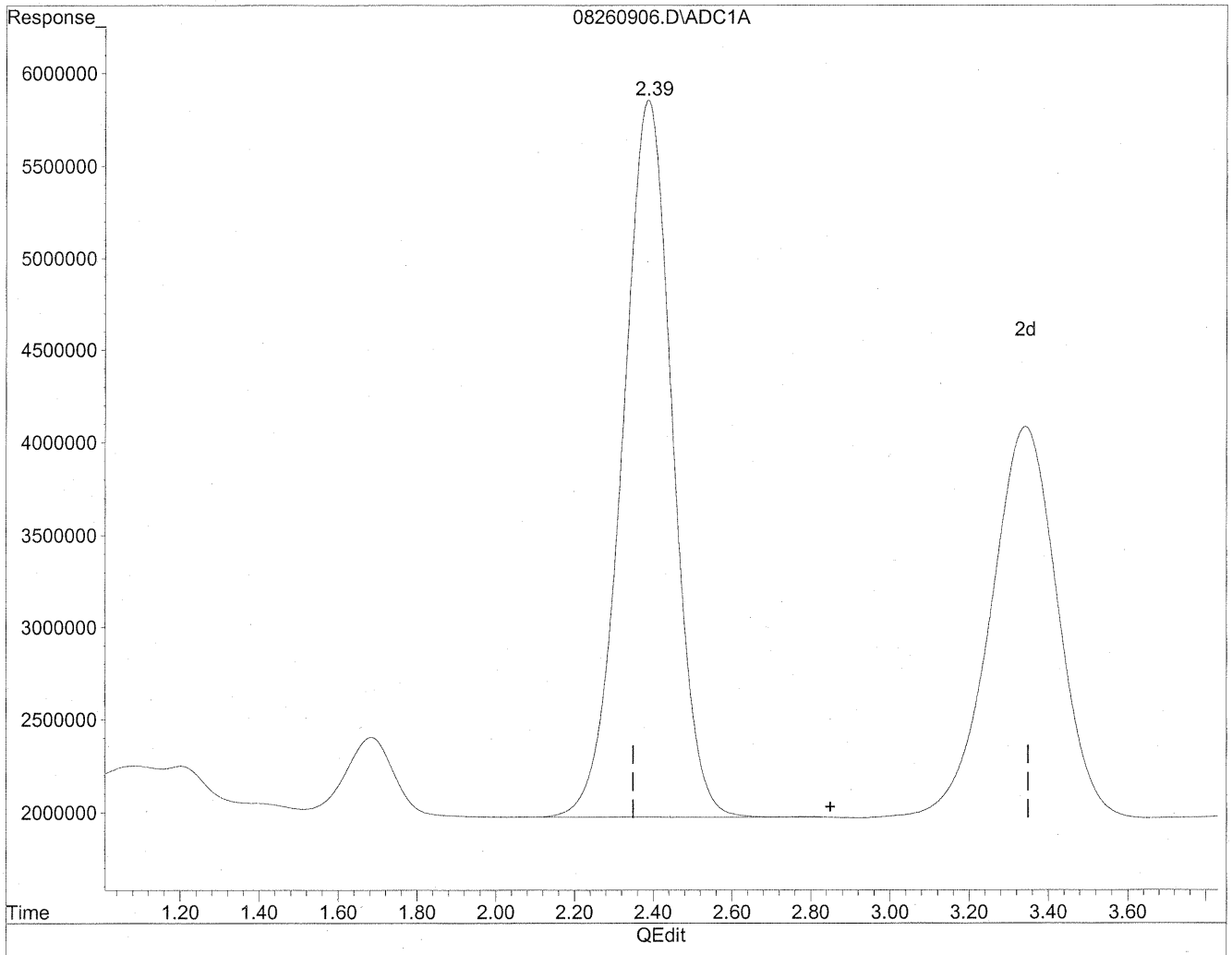
8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260906.D Vial: 6
Acq On : 26 Aug 2009 6:20 pm Operator: HC
Sample : P0902942-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:45 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.39min 3210.775ng/ml

response 342574387

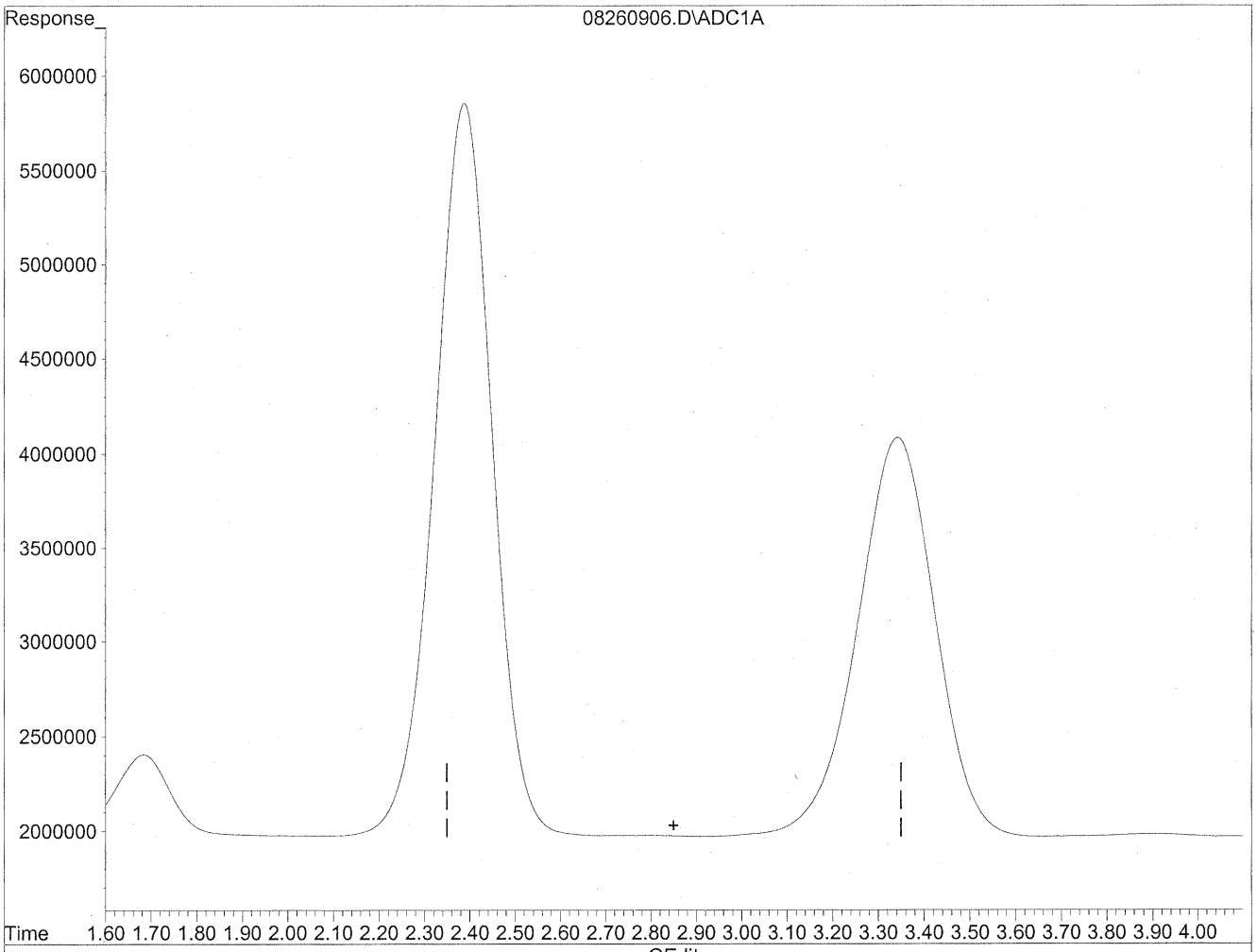
(+) = Expected Retention Time

08260906.D TO110709.M Sat Aug 29 16:46:24 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260906.D Vial: 6
Acq On : 26 Aug 2009 6:20 pm Operator: HC
Sample : P0902942-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:45 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/30/09

LA 8/31/09

33

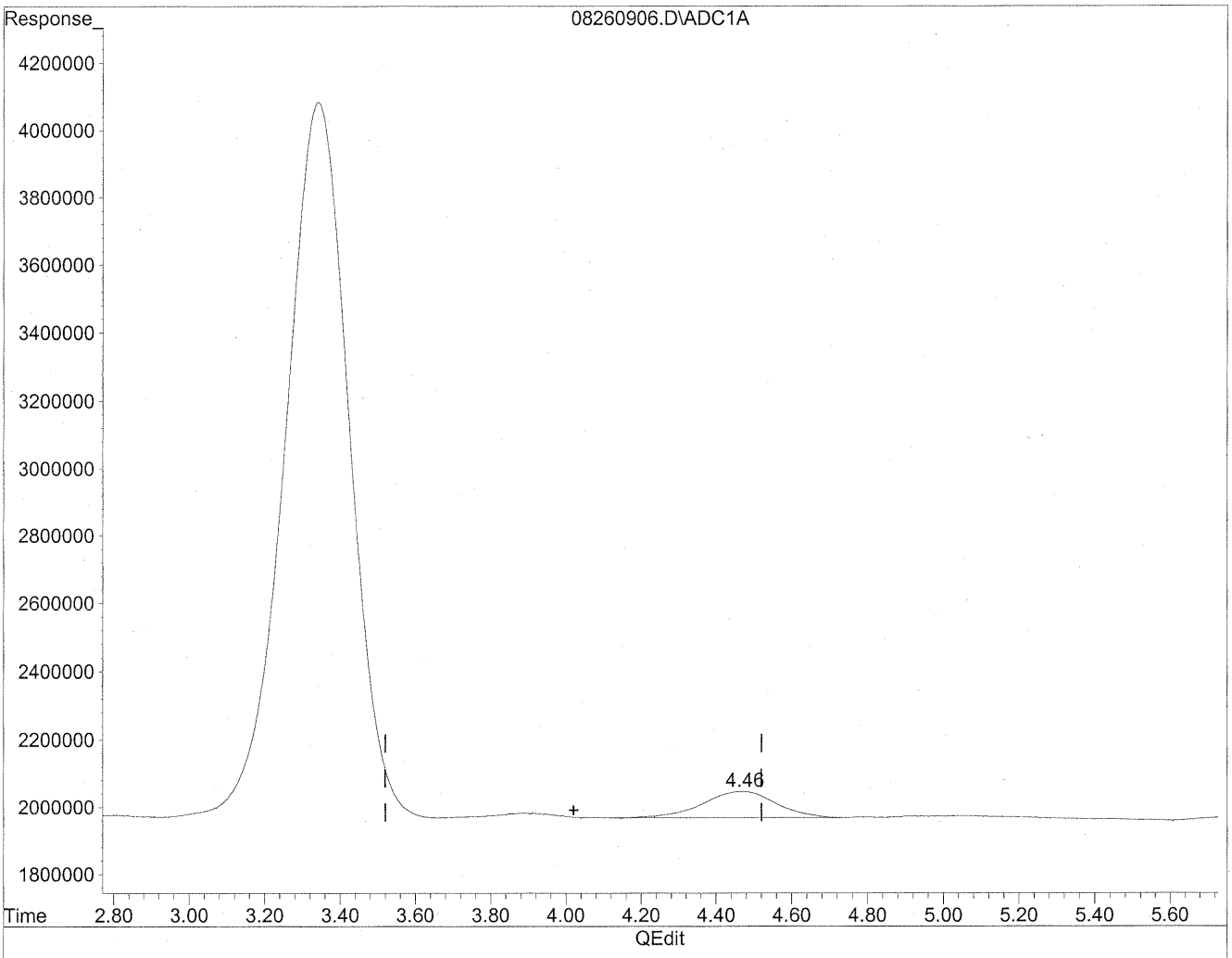
(+) = Expected Retention Time

08260906.D TO110709.M Sat Aug 29 16:46:28 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260906.D Vial: 6
Acq On : 26 Aug 2009 6:20 pm Operator: HC
Sample : P0902942-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:45 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.47min 107.481ng/ml

response 10470289

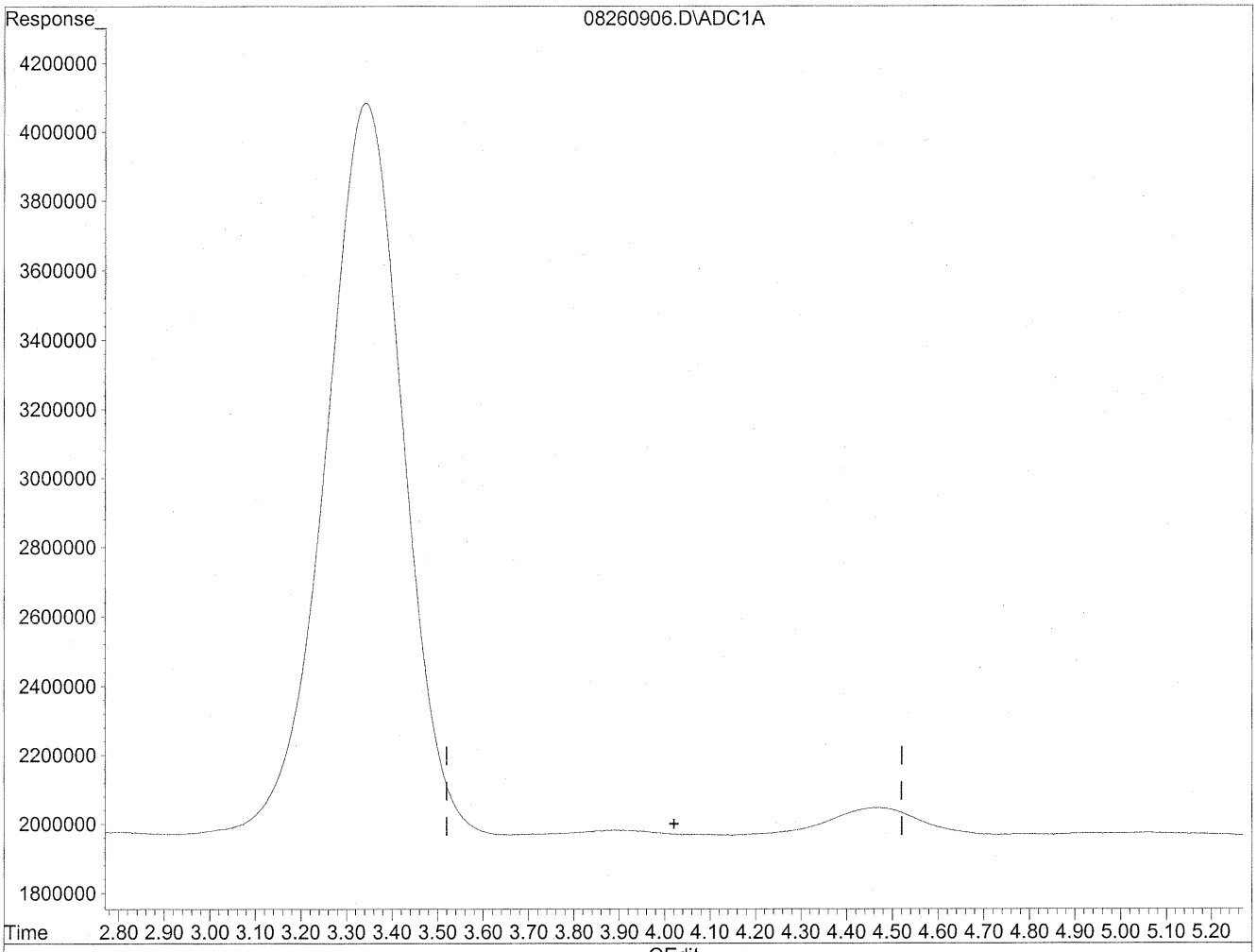
(+) = Expected Retention Time

08260906.D TO110709.M Sat Aug 29 16:46:32 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260906.D Vial: 6
Acq On : 26 Aug 2009 6:20 pm Operator: HC
Sample : P0902942-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:45 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

tlc 8/30/09 wp

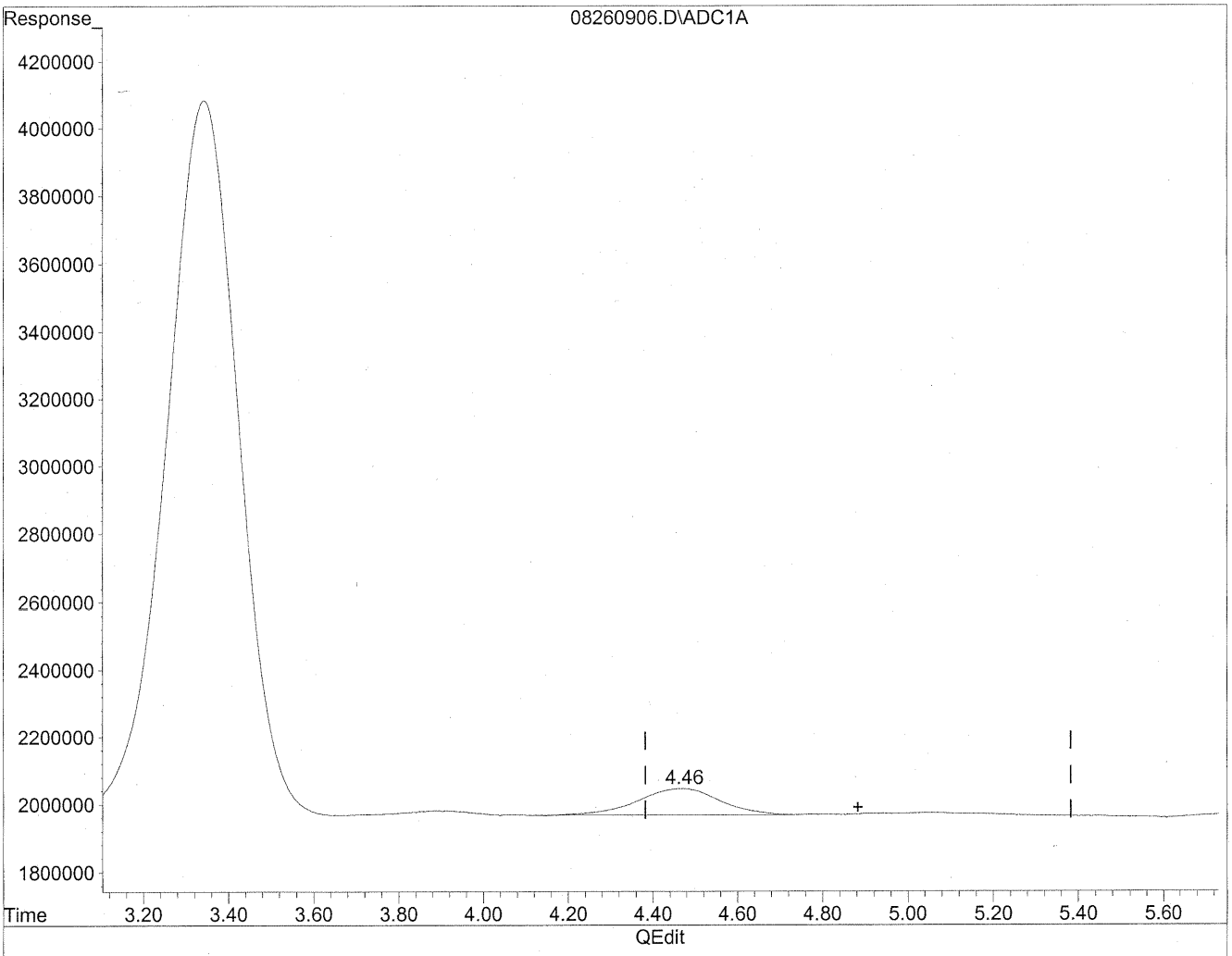
8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260906.D Vial: 6
Acq On : 26 Aug 2009 6:20 pm Operator: HC
Sample : P0902942-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:45 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.47min 118.528ng/ml

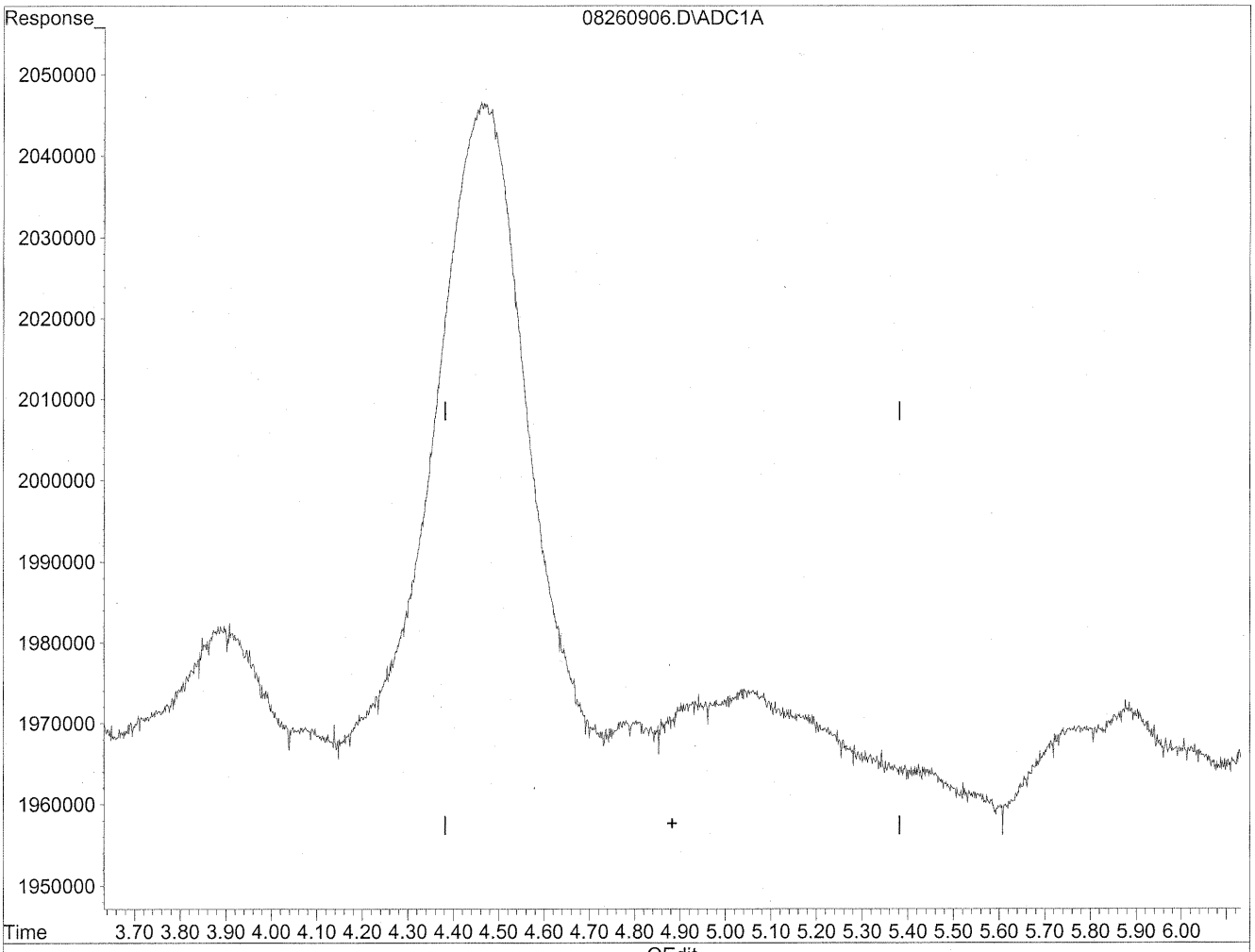
response 10470289

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260906.D Vial: 6
Acq On : 26 Aug 2009 6:20 pm Operator: HC
Sample : P0902942-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:45 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



QEdit

(5) Butyraldehyde

0.00min 0.000ng/ml d

response 0

HC 8/30/09 MP

8/31/09

(+) = Expected Retention Time

08260906.D TO110709.M Sat Aug 29 16:46:42 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **101470**
Client Project ID: **16512**

CAS Project ID: P0902942
CAS Sample ID: P0902942-002

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: 8/24/09
Date Received: 8/25/09
Date Analyzed: 8/26-27/09
Desorption Volume: 1.0 ml
Volume Sampled: 104.04 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**2,300**	**22**	0.96	**18**	0.78	
75-07-0	**Acetaldehyde**	**4,700**	**46**	0.96	**25**	0.53	
123-38-6	**Propionaldehyde**	**570**	**5.5**	0.96	**2.3**	0.40	
4170-30-3	Crotonaldehyde, Total	< 200	ND	1.9	ND	0.67	**V**
123-72-8	**Butyraldehyde**	**1,100**	**10**	0.96	**3.5**	0.33	**M**
100-52-7	**Benzaldehyde**	**840**	**8.1**	0.96	**1.9**	0.22	
590-86-3	**Isovaleraldehyde**	**170**	**1.6**	0.96	**0.47**	0.27	
110-62-3	**Valeraldehyde**	**1,500**	**15**	0.96	**4.2**	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**	**4,600**	**44**	0.96	**11**	0.23	
5779-94-2	2,5-Dimethylbenzaldehyde	< 200	ND	1.9	ND	0.35	**V**

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.
M = Matrix interference; results may be biased high.
V = The continuing calibration verification standard was outside (biased low) the specified limits for this compound.

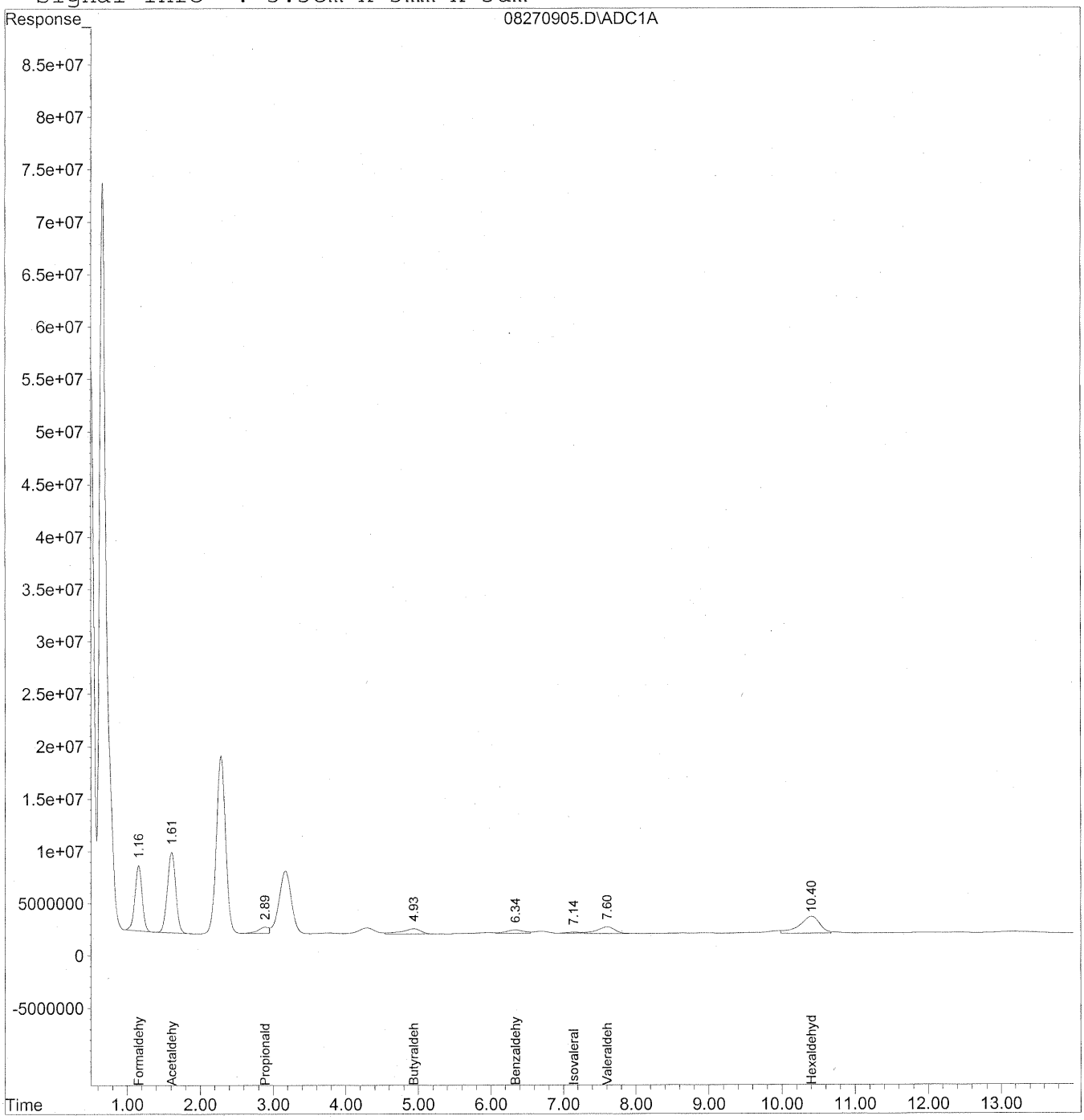
Verified By: RG Date: 9/11/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:15 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 : Thu Aug 27 08:33:51 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\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:15 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 : Thu Aug 27 08:33:51 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

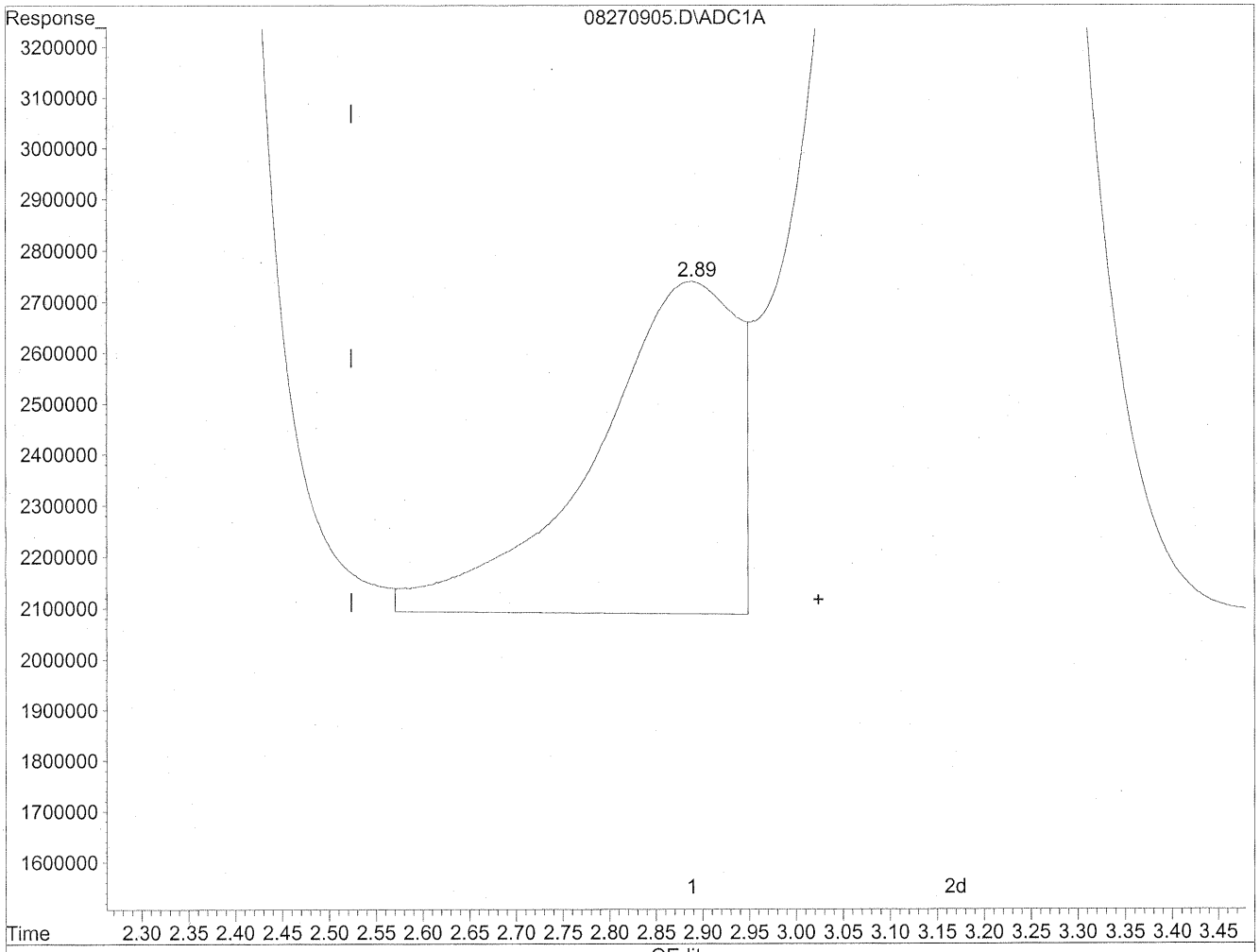
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	426986739	2325.872	ng/ml
2)	Acetaldehyde	1.61	623235677	4444.588	ng/ml
3)	Propionaldehyde	2.89	60514643	567.173	ng/mlm
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	4.93	94605293	1070.970	ng/mlm
6)	Benzaldehyde	6.34	55299022	839.526	ng/mlm
7)	Isovaleraldehyde	7.14	13432794	171.663	ng/mlm
8)	Valeraldehyde	7.60	112933756	1536.409	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	10.40f	308320168	4578.303	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\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.89min 651.607ng/ml

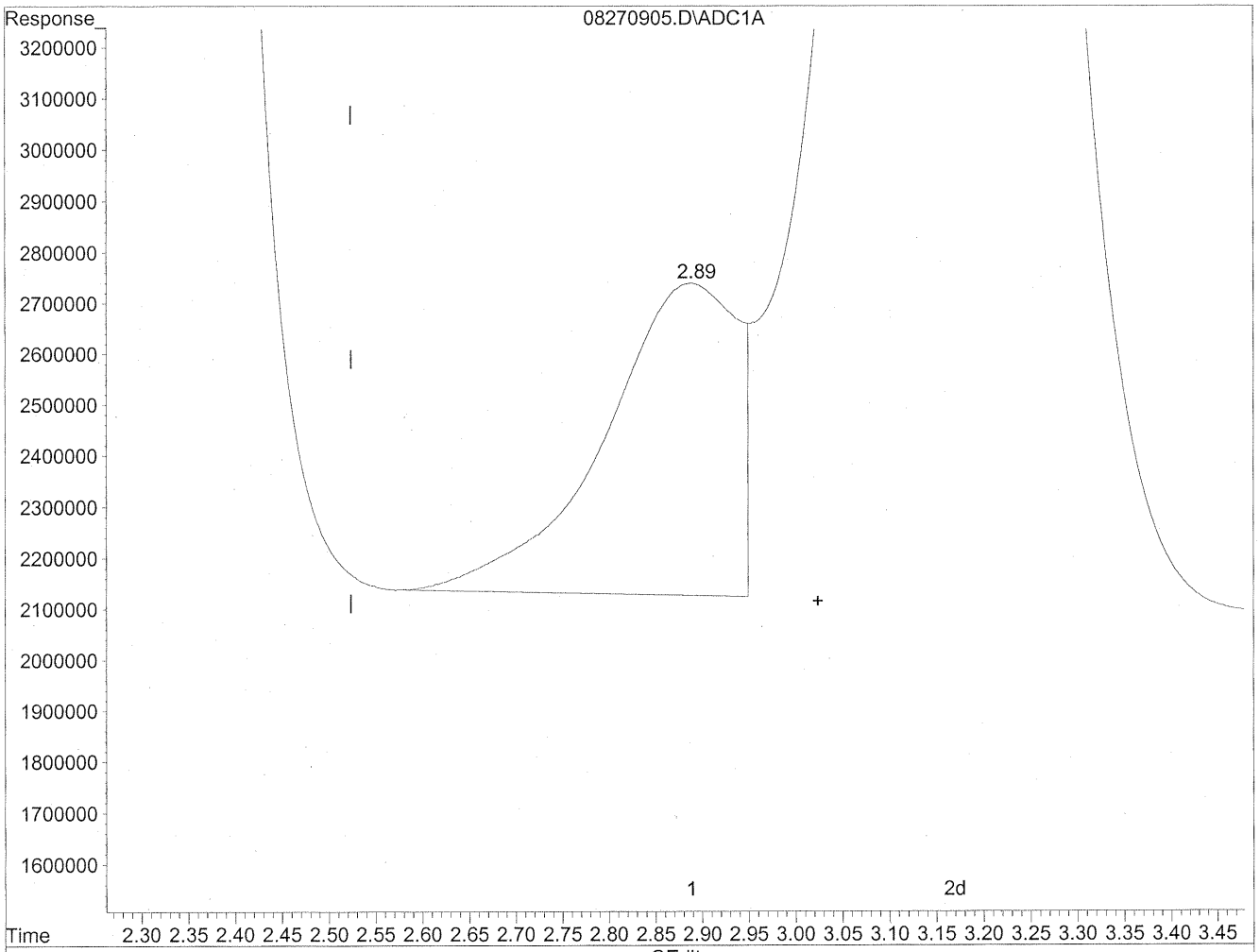
response 69523321

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.89min 567.173ng/ml m

response 60514643

HC 8/31/09

VK

WA 8/31/09

42

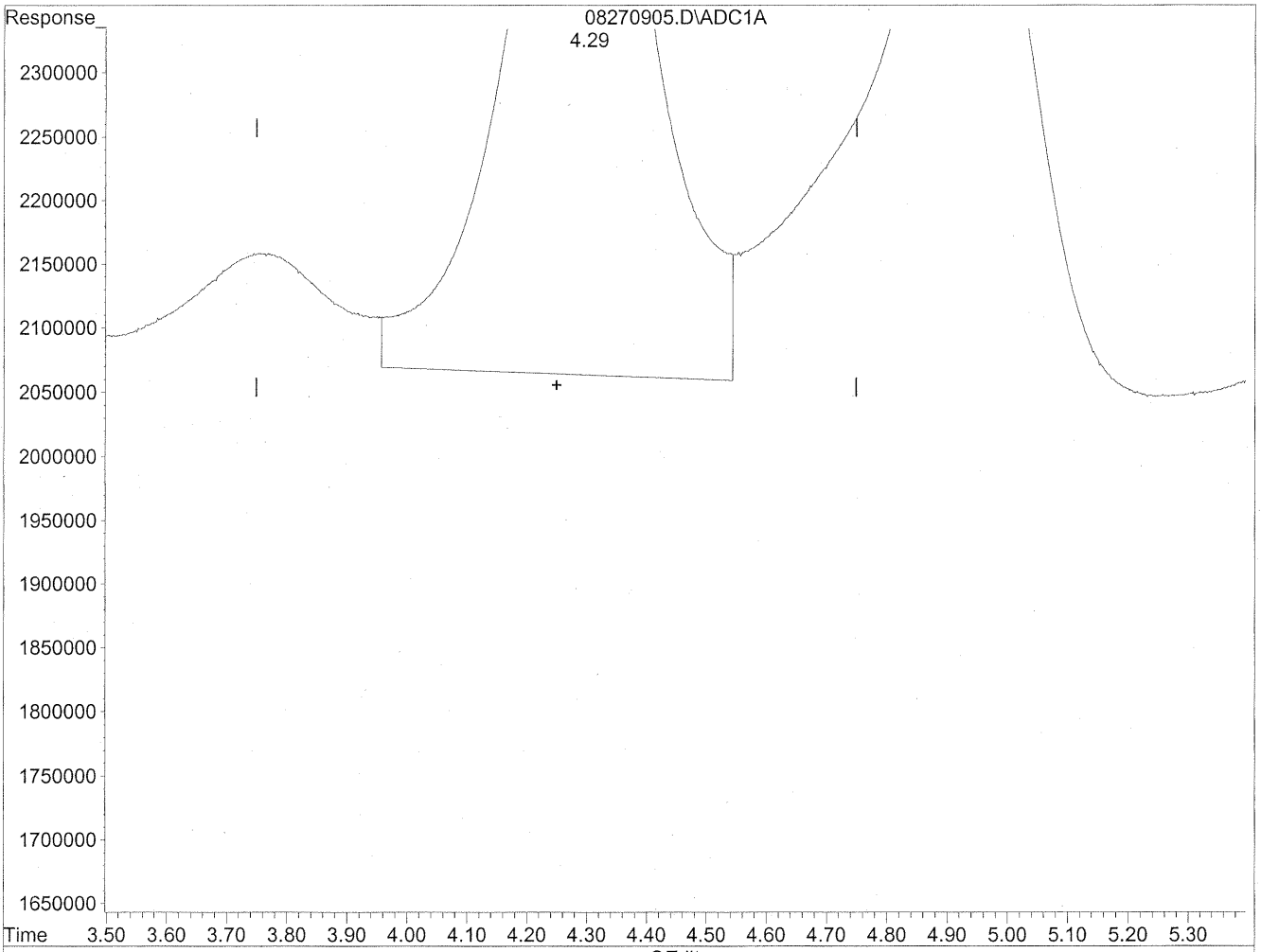
(+) = Expected Retention Time

08270905.D TO110709.M Sun Aug 30 13:12:32 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.29min 951.983ng/ml

response 92737634

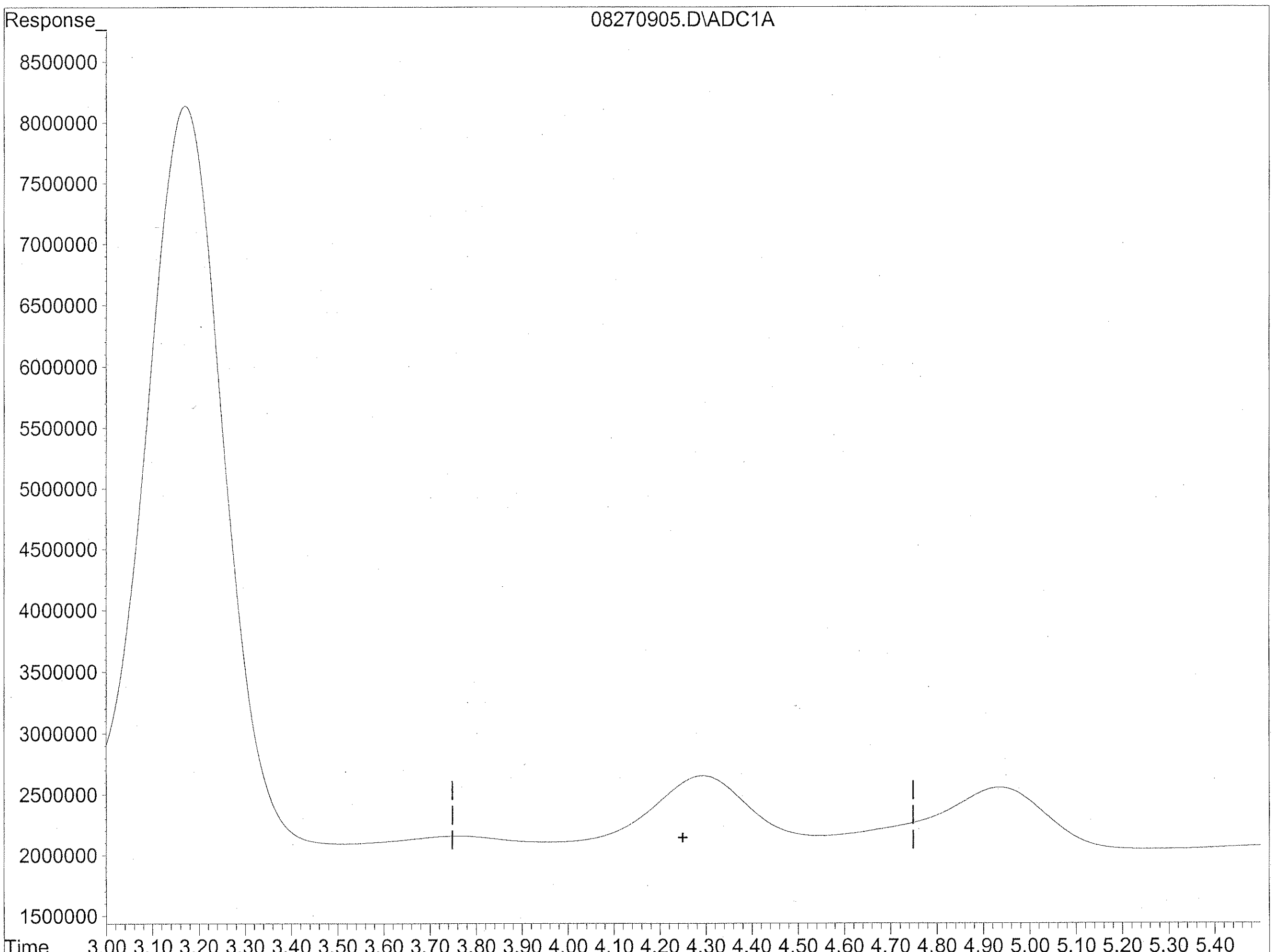
(+) = Expected Retention Time

08270905.D TO110709.M Sun Aug 30 13:12:36 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



QEdit

(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/31/09

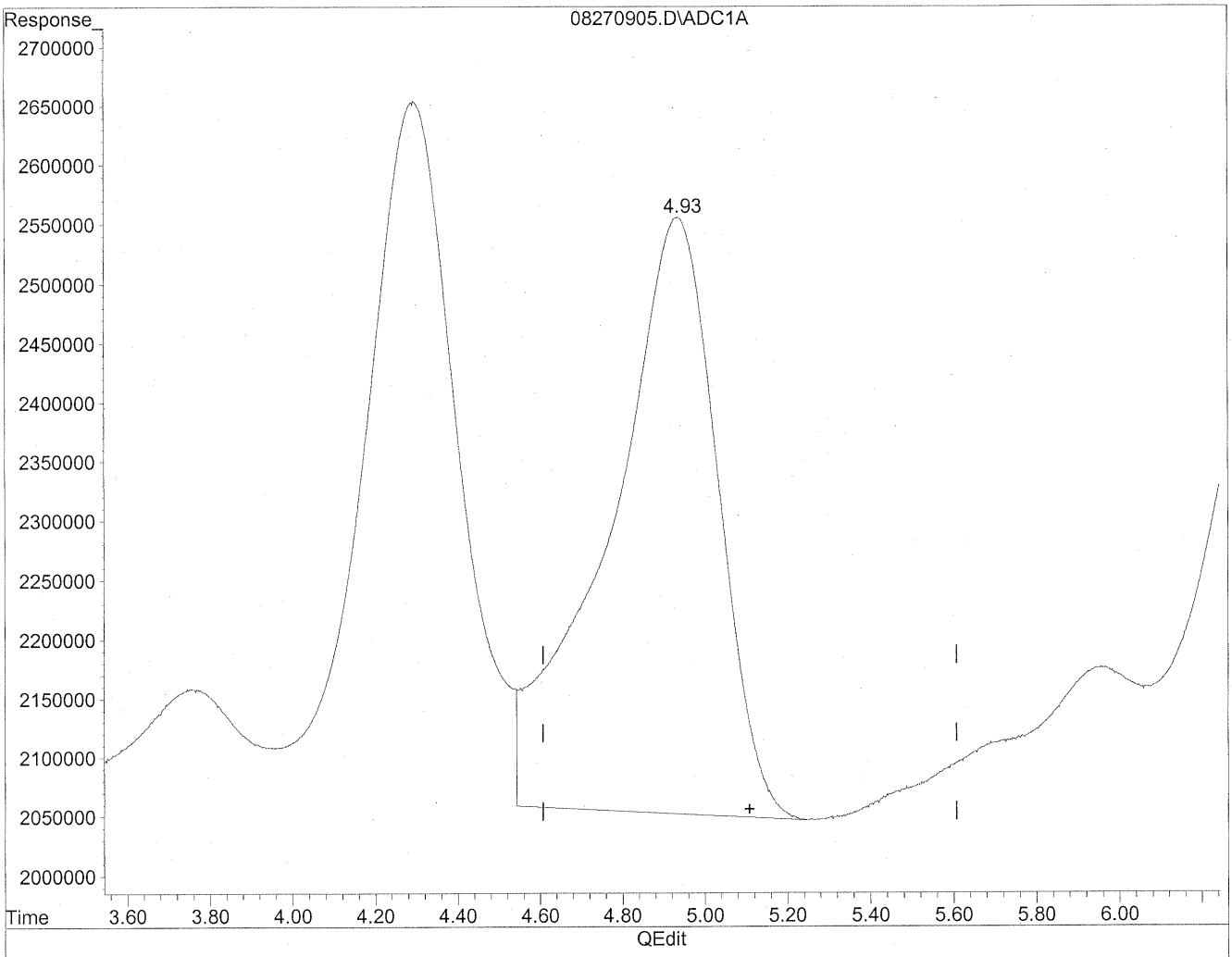
(+) = Expected Retention Time

08270905.D TO110709.M Sun Aug 30 13:12:38 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.93min 1040.766ng/ml

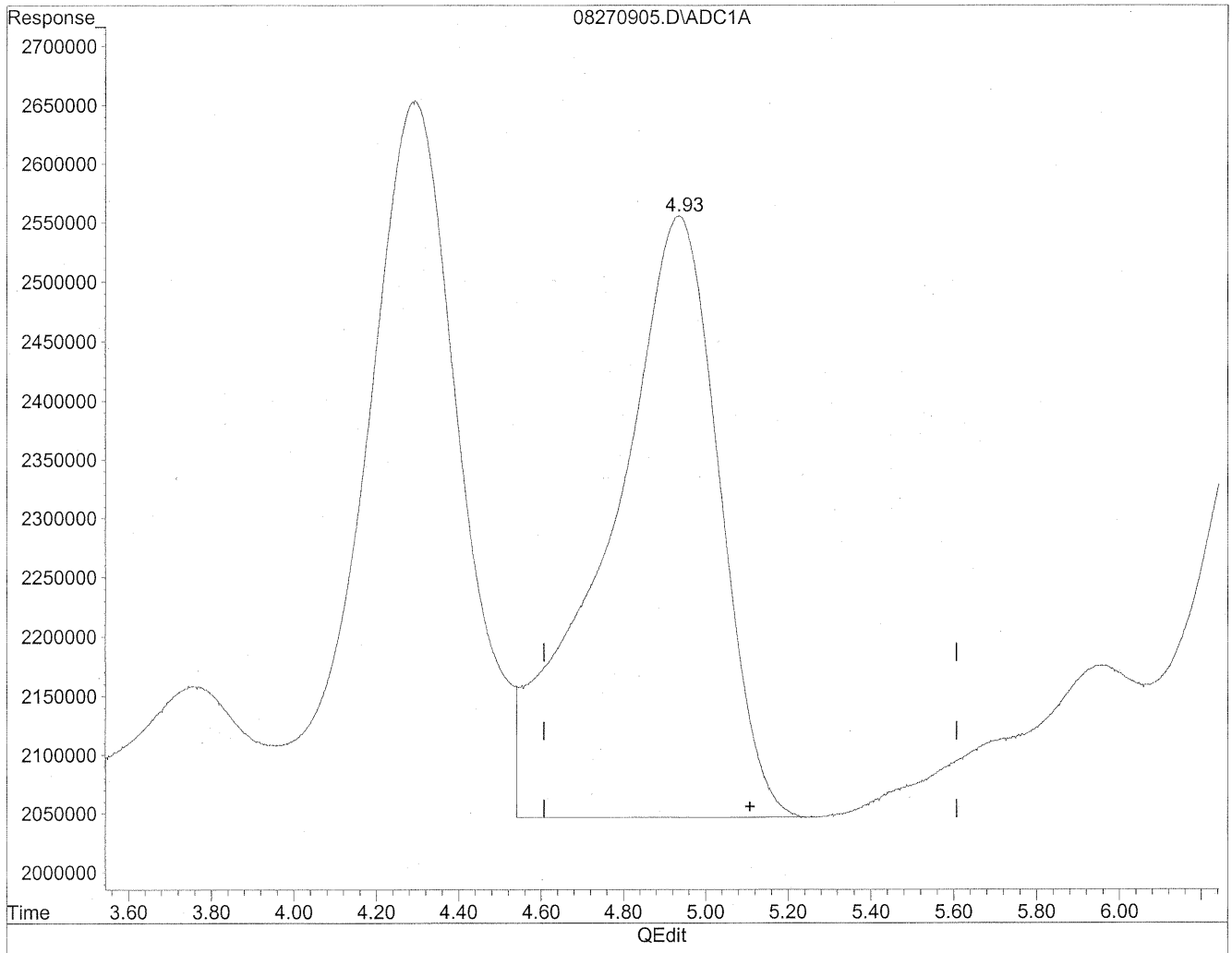
response 91937201

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.93min 1070.970ng/ml m

response 94605293

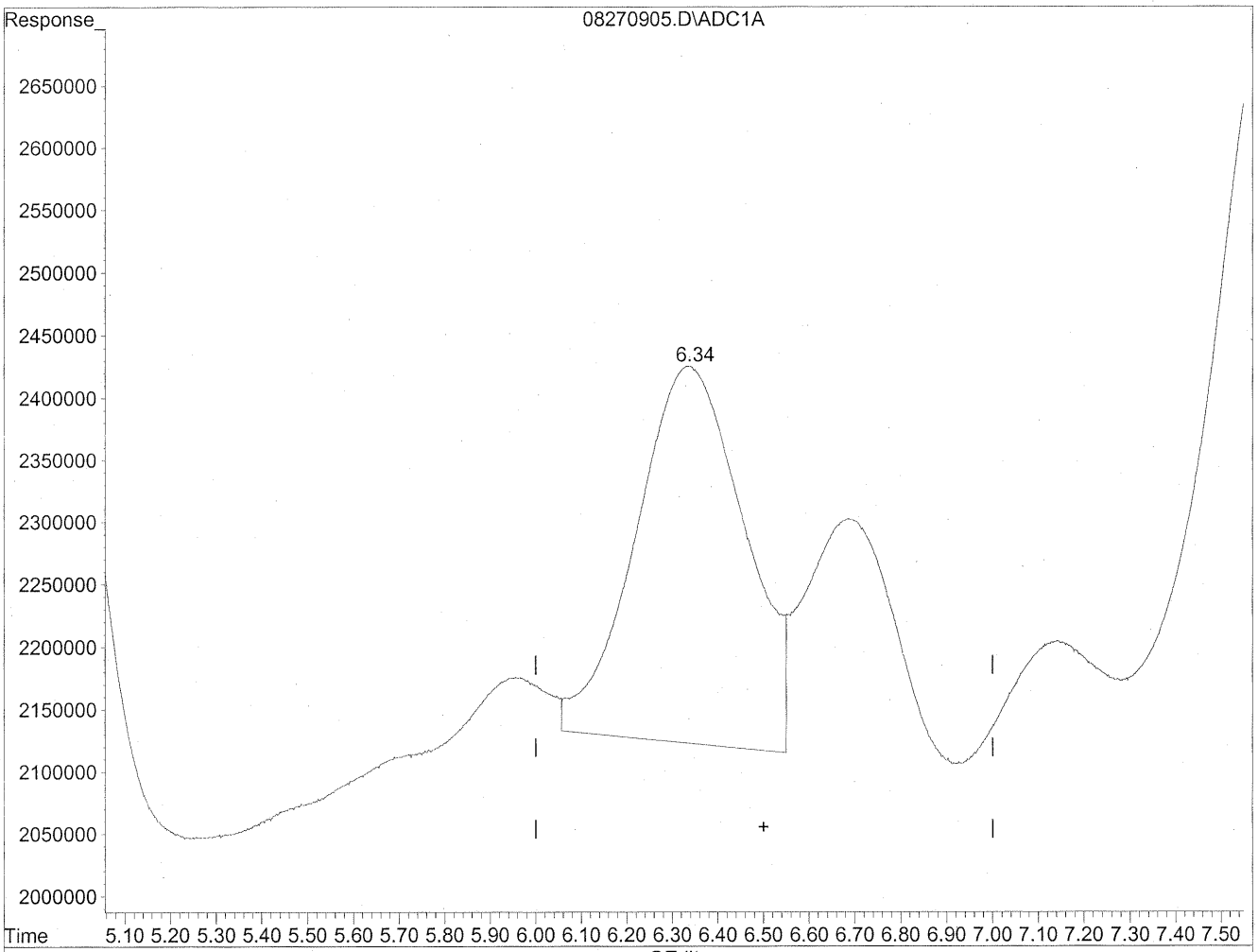
HC 8/31/09
BC
MP
8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.34min 761.018ng/ml

response 50127726

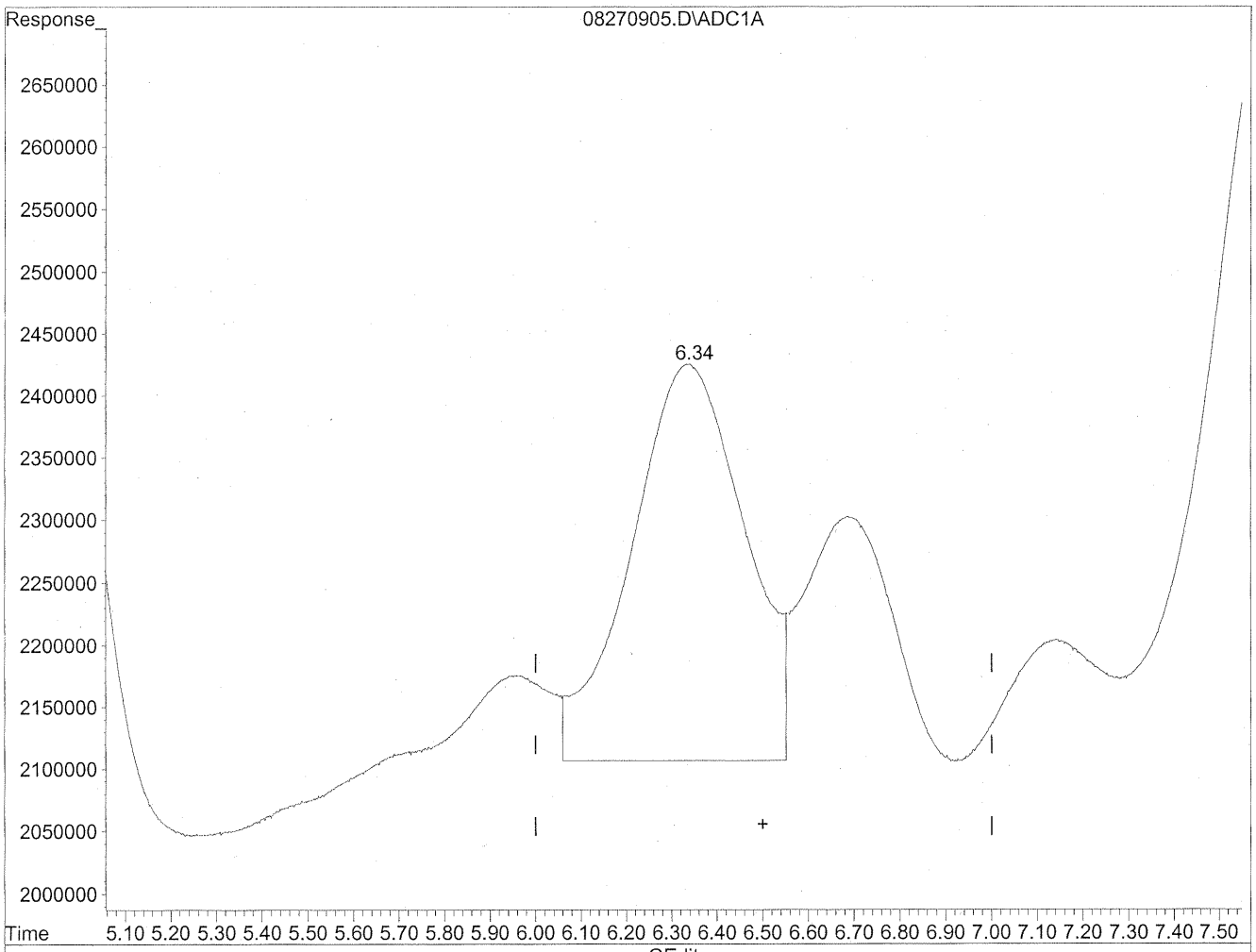
(+) = Expected Retention Time

08270905.D TO110709.M Sun Aug 30 13:12:59 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.34min 839.526ng/ml m

response 55299022

HC 8/31/09
BC

8/31/09

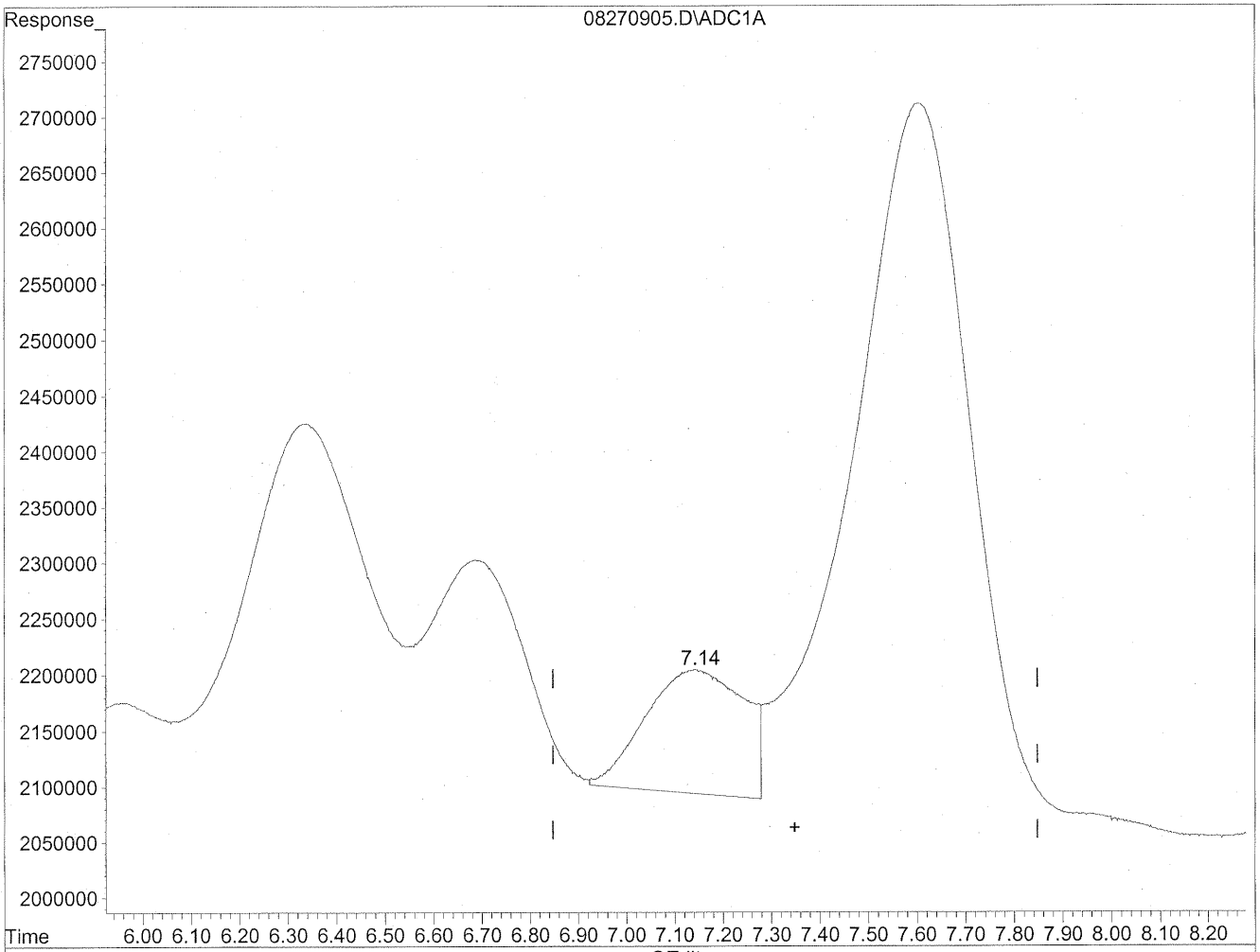
48

(+) = Expected Retention Time
08270905.D TO110709.M Sun Aug 30 13:13:32 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.14min 200.265ng/ml

response 15670906

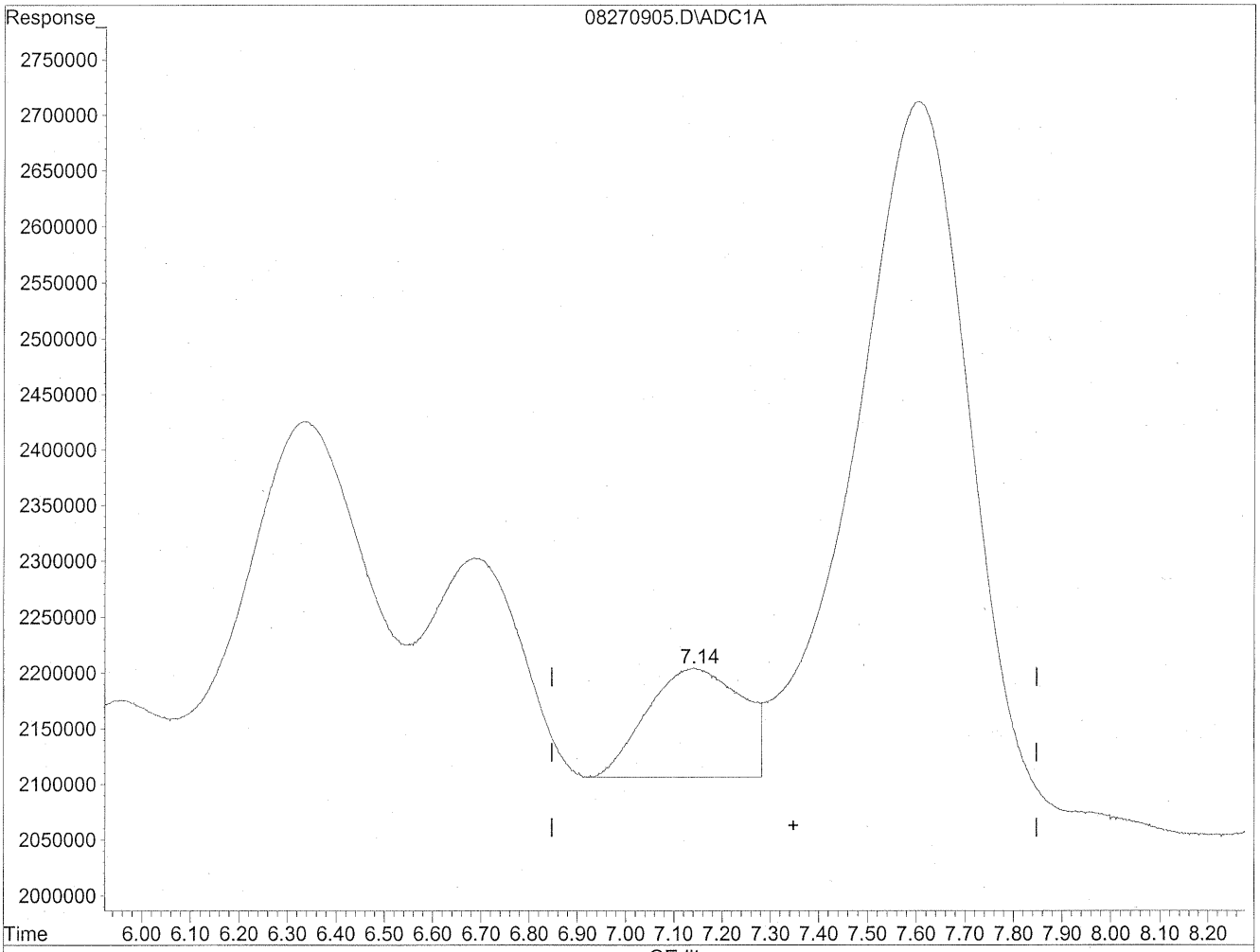
(+) = Expected Retention Time

08270905.D TO110709.M Sun Aug 30 13:13:36 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.14min 171.663ng/ml m

response 13432794

T/C 8/31/09 IC

8/31/09

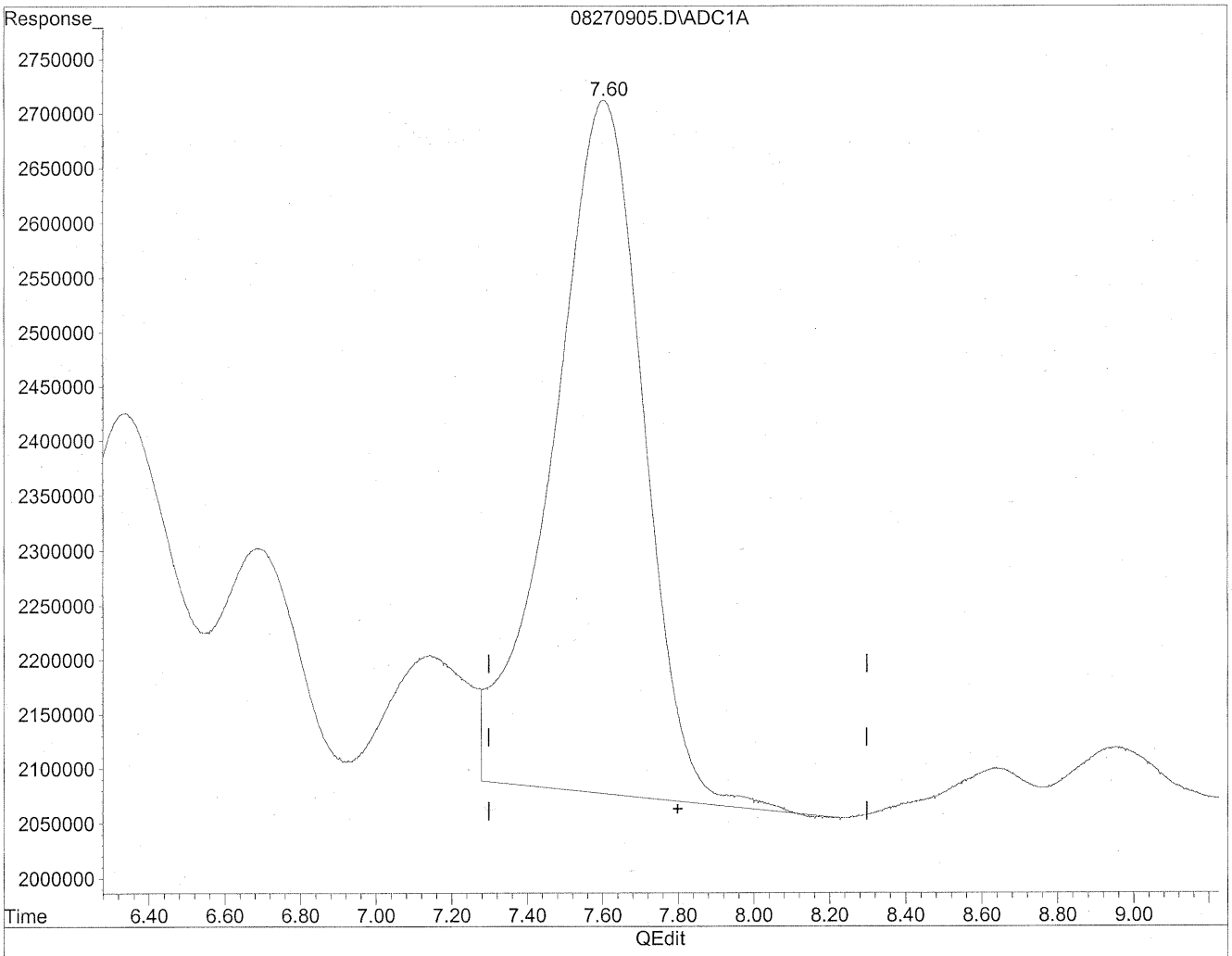
(+) = Expected Retention Time

08270905.D TO110709.M Sun Aug 30 13:13:48 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.60min 1444.434ng/ml

response 106173113

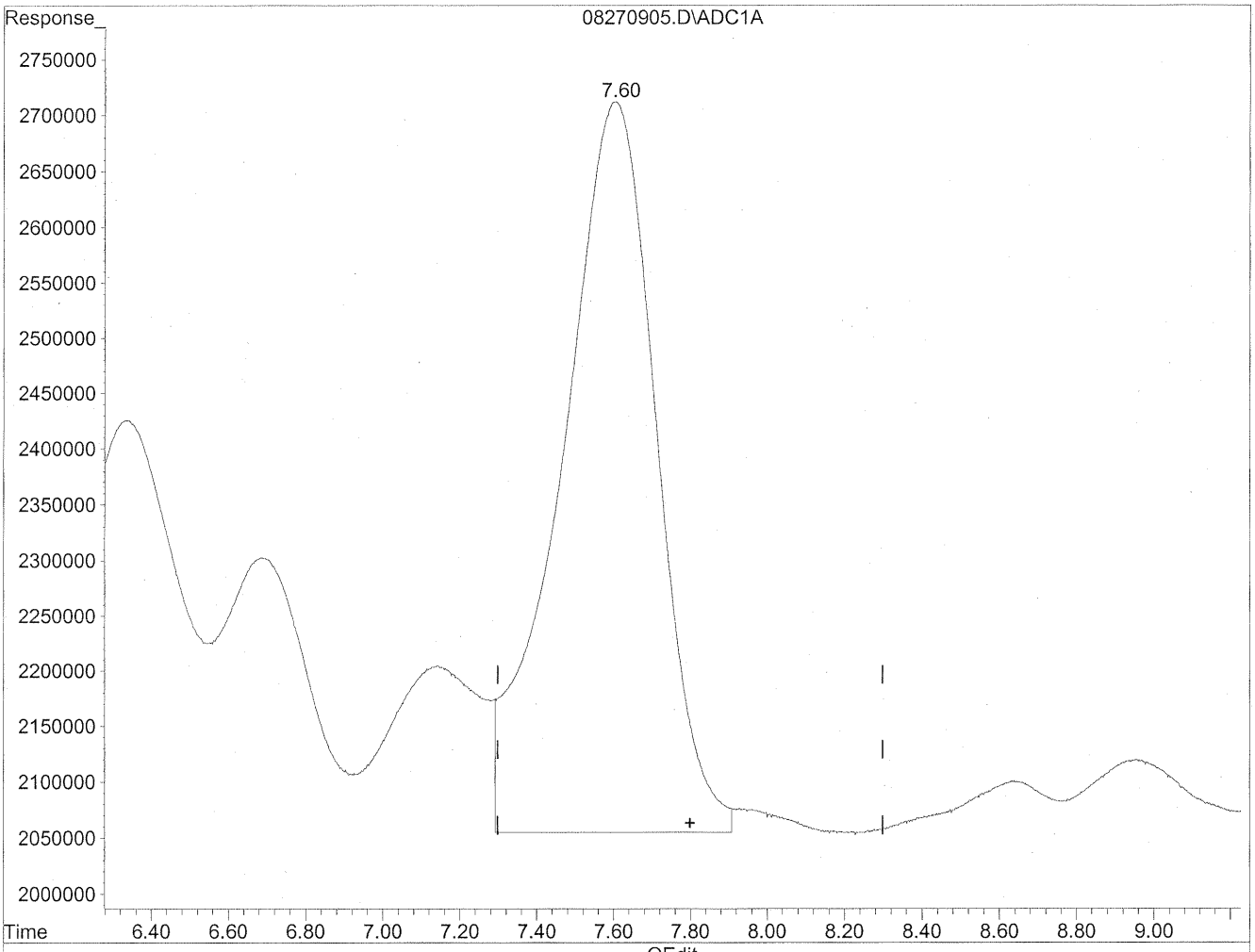
(+) = Expected Retention Time

08270905.D TO110709.M Sun Aug 30 13:13:51 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.60min 1536.409ng/ml m

response 112933756

HC 8/31/09
SH, BC

8/31/09

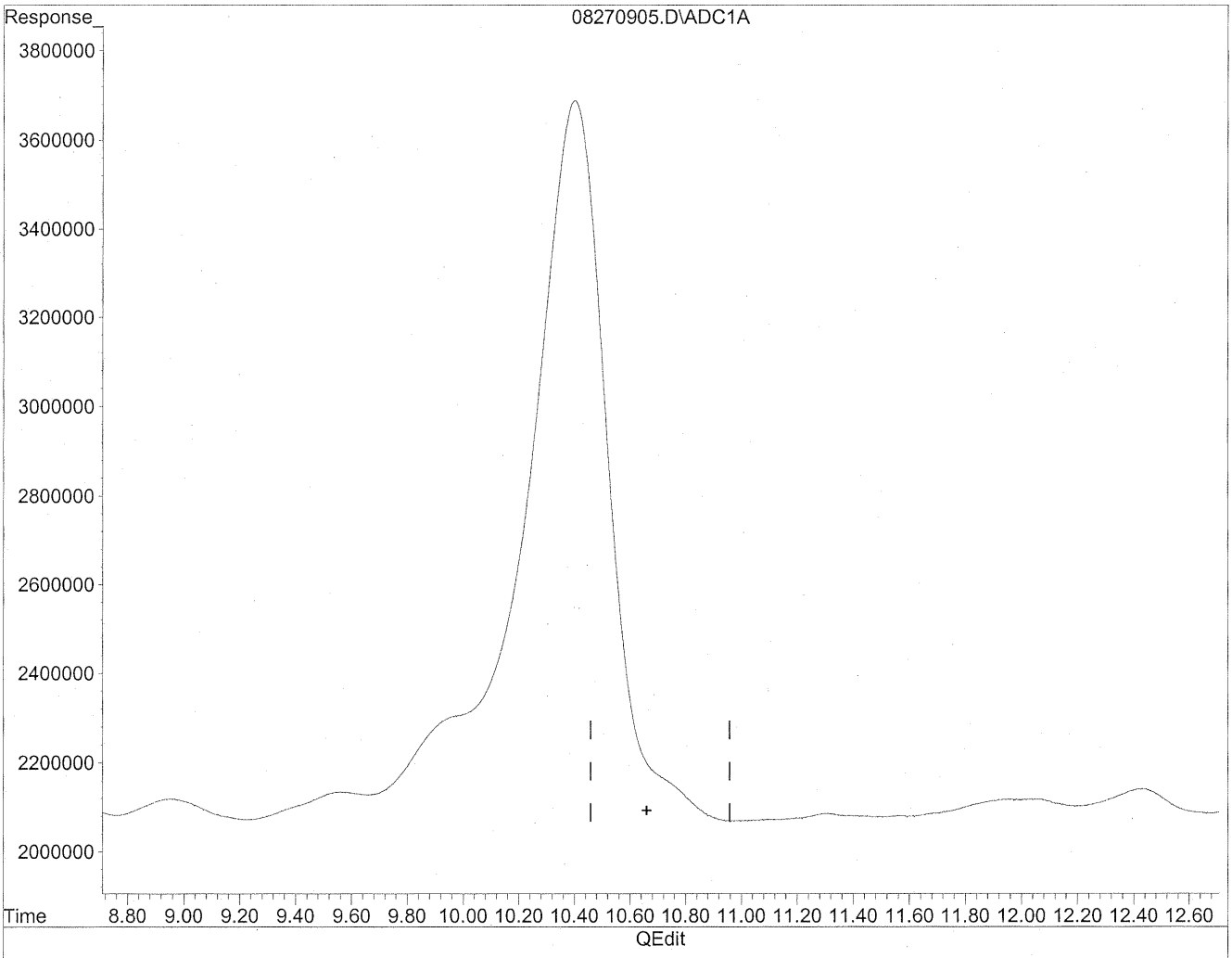
(+) = Expected Retention Time

08270905.D TO110709.M Sun Aug 30 13:14:16 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.66min 0.000ng/ml

response 0

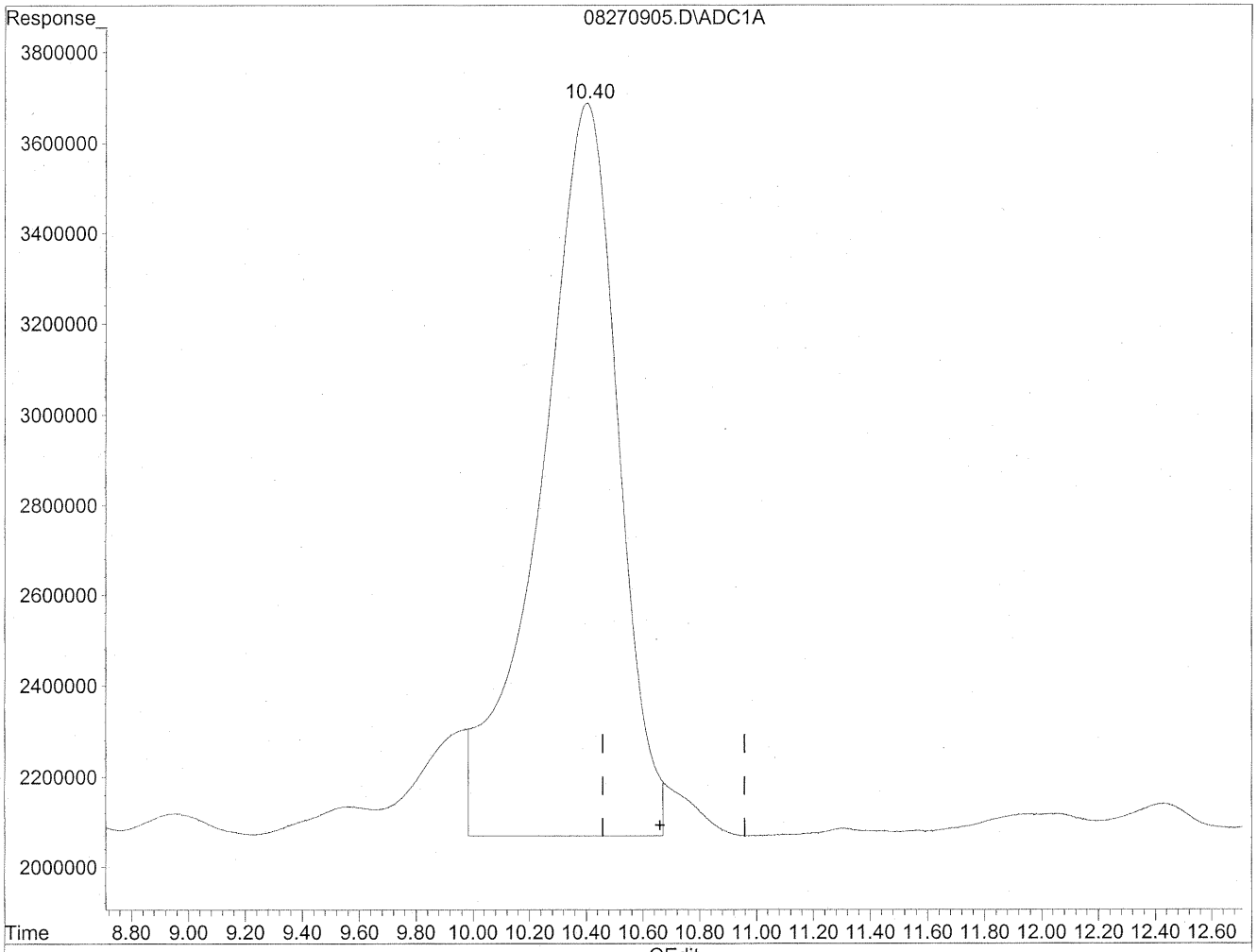
(+) = Expected Retention Time

08270905.D TO110709.M Sun Aug 30 13:14:23 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270905.D Vial: 5
Acq On : 27 Aug 2009 10:05 am Operator: HC
Sample : P0902942-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 10:25 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.40min 4578.303ng/ml m

response 308320168

tlc 8/31/09 BN

8/31/09

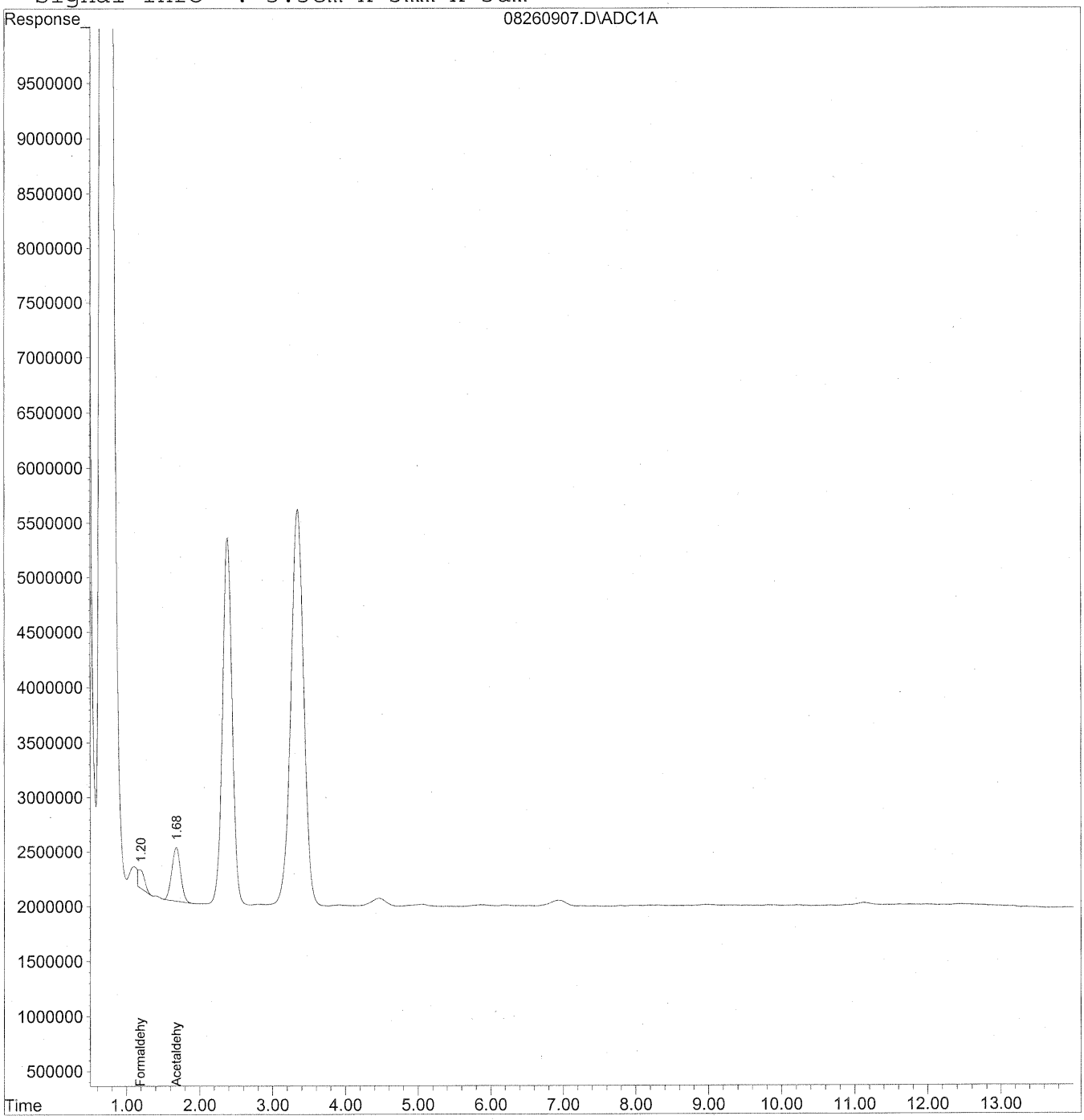
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260907.D Vial: 7
Acq On : 26 Aug 2009 6:35 pm Operator: HC
Sample : P0902942-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:47 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 : Thu Aug 27 07:35:56 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



55

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\26\08260907.D Vial: 7
Acq On : 26 Aug 2009 6:35 pm Operator: HC
Sample : P0902942-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:47 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 : Thu Aug 27 07:35:56 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

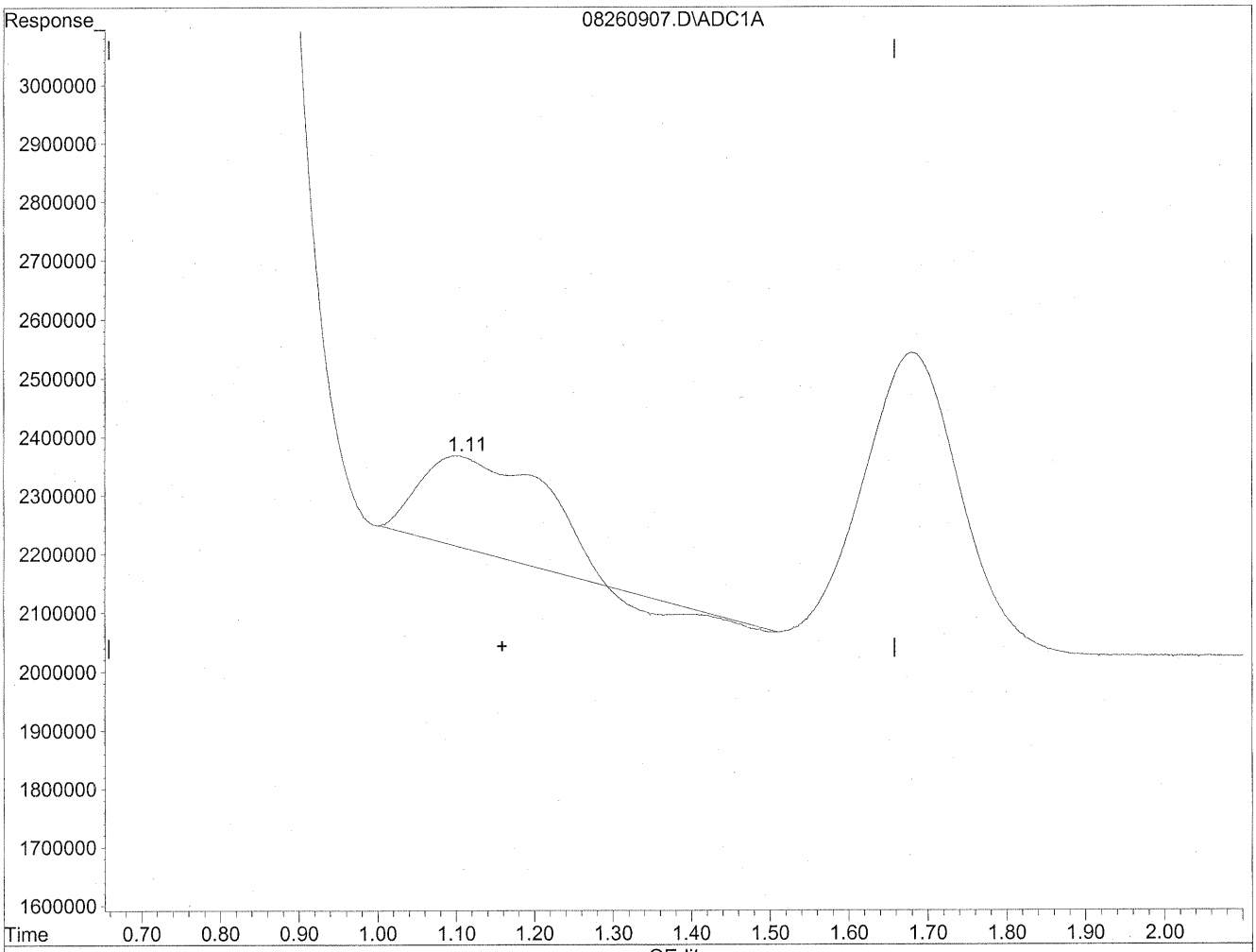
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.20	10438462	56.860	ng/mlm
2)	Acetaldehyde	1.68	41522942	296.120	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/mld
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\26\08260907.D Vial: 7
Acq On : 26 Aug 2009 6:35 pm Operator: HC
Sample : P0902942-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.10min 93.466ng/ml

response 17158529

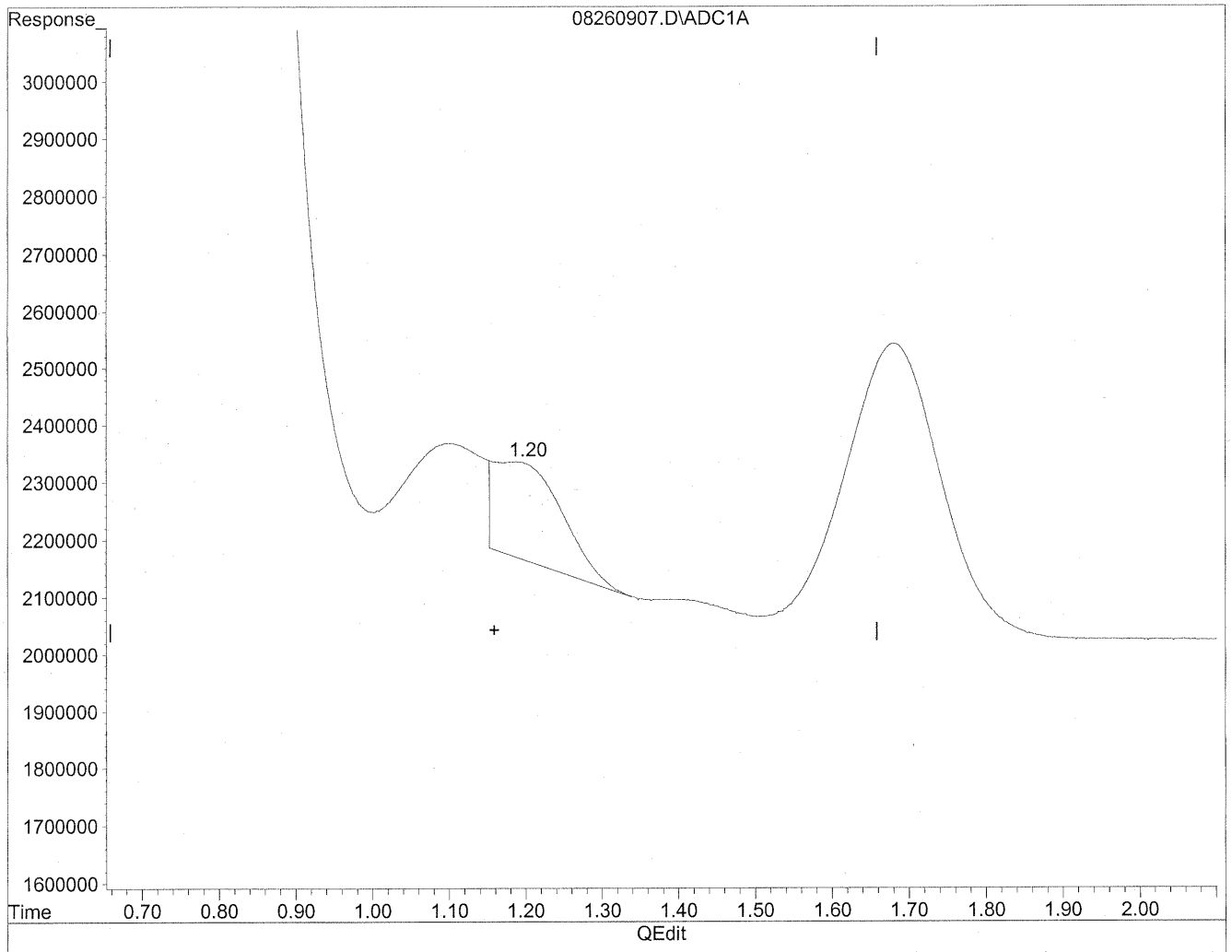
(+) = Expected Retention Time

08260907.D TO110709.M Sat Aug 29 16:47:12 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260907.D Vial: 7
Acq On : 26 Aug 2009 6:35 pm Operator: HC
Sample : P0902942-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.20min 56.860ng/ml m

response 10438462

HC 8/30/09

SP

MB 8/31/09

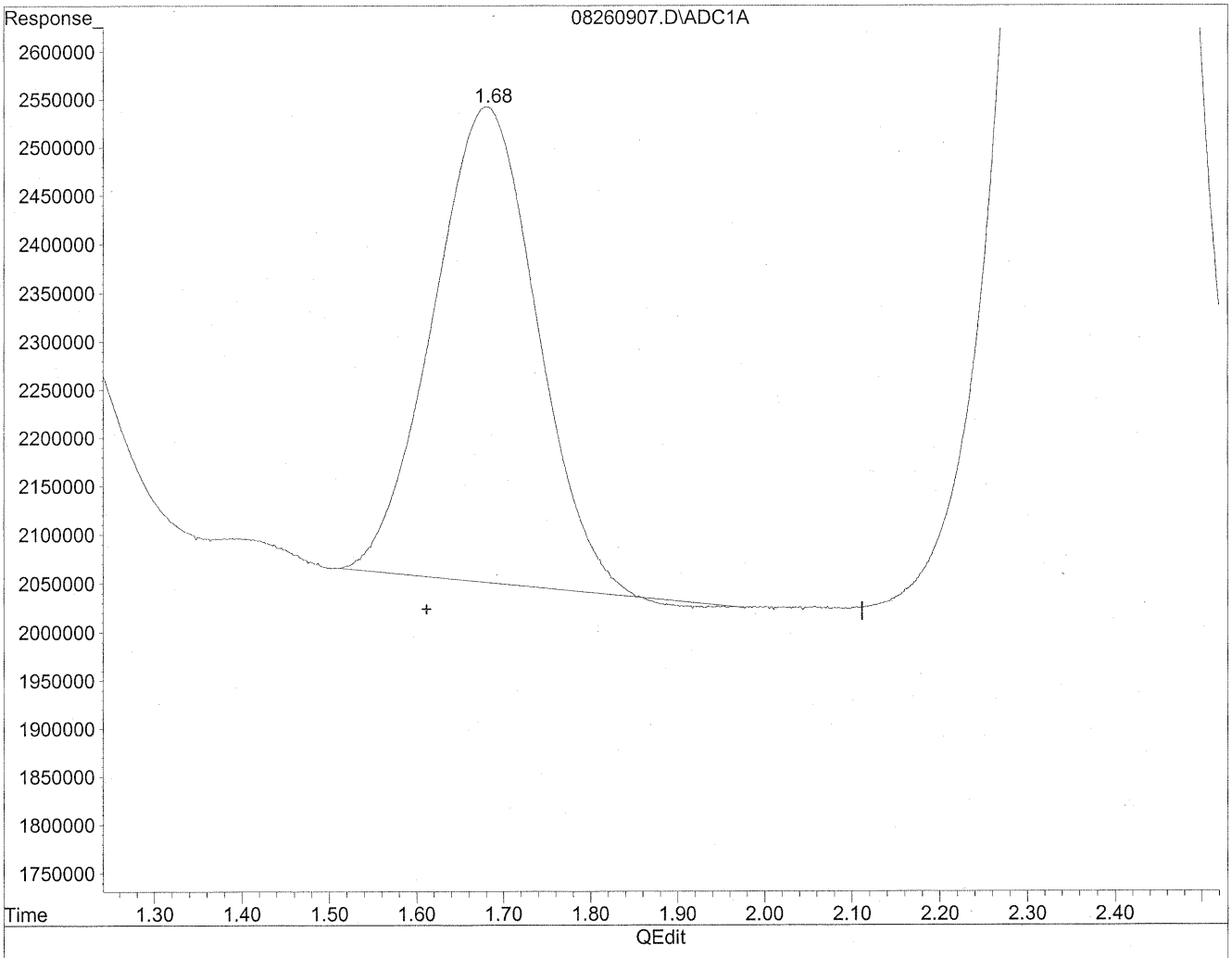
(+) = Expected Retention Time

08260907.D TO110709.M Sat Aug 29 16:47:19 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260907.D Vial: 7
Acq On : 26 Aug 2009 6:35 pm Operator: HC
Sample : P0902942-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.68min 290.191ng/ml

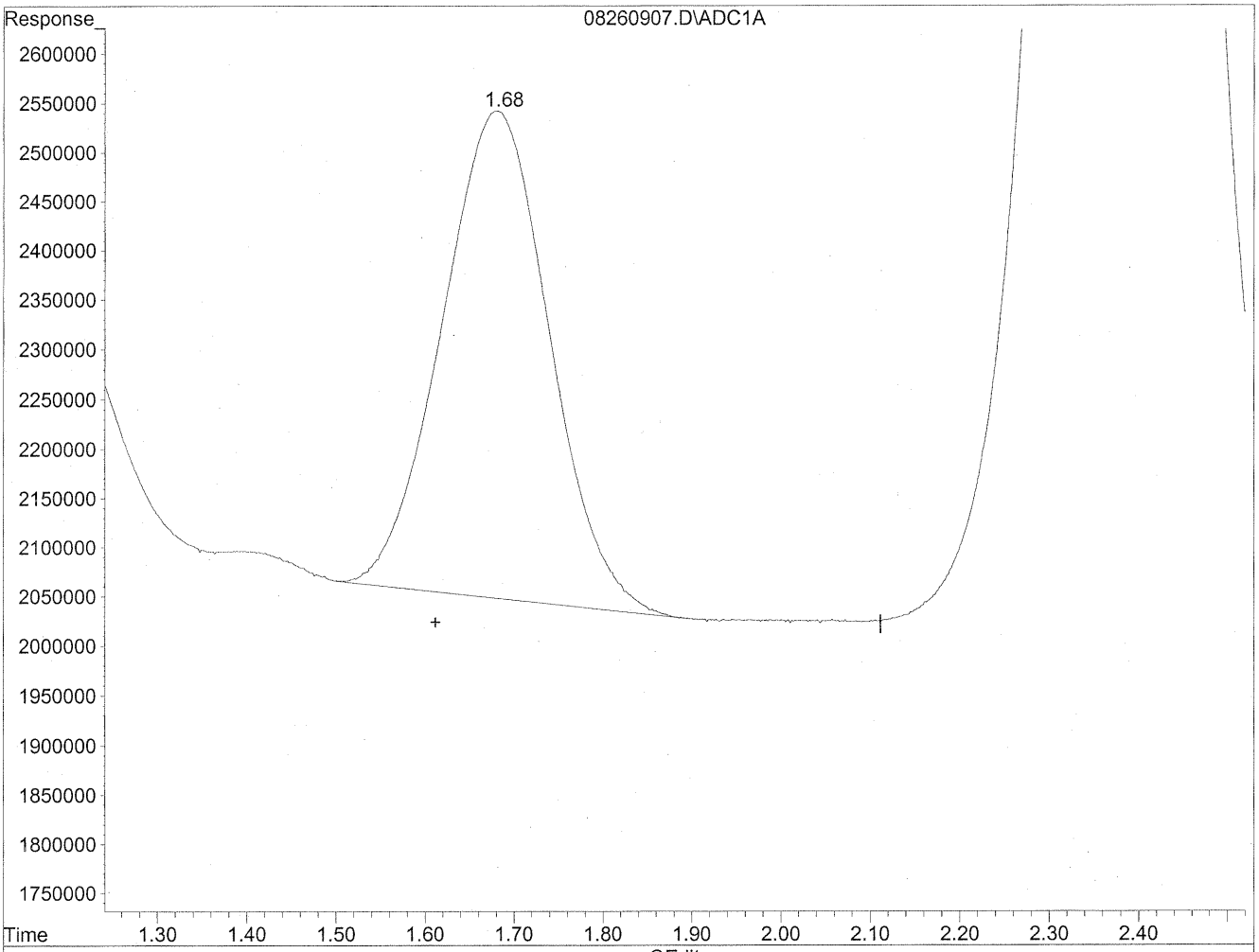
response 40691560

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260907.D Vial: 7
Acq On : 26 Aug 2009 6:35 pm Operator: HC
Sample : P0902942-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.68min 296.120ng/ml m

response 41522942

HC 8/30/09

LC

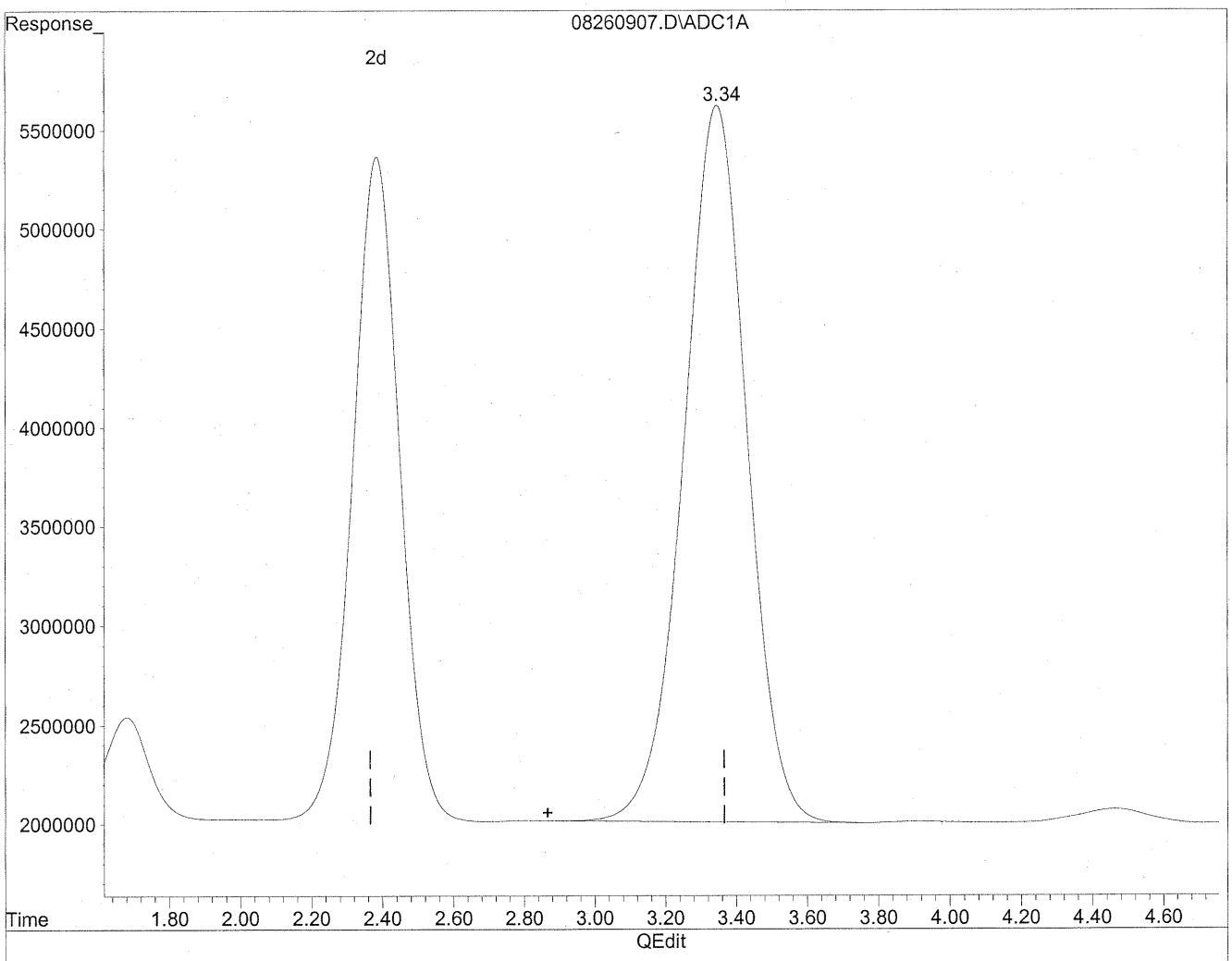
CW 8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260907.D Vial: 7
Acq On : 26 Aug 2009 6:35 pm Operator: HC
Sample : P0902942-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.35min 4187.069ng/ml

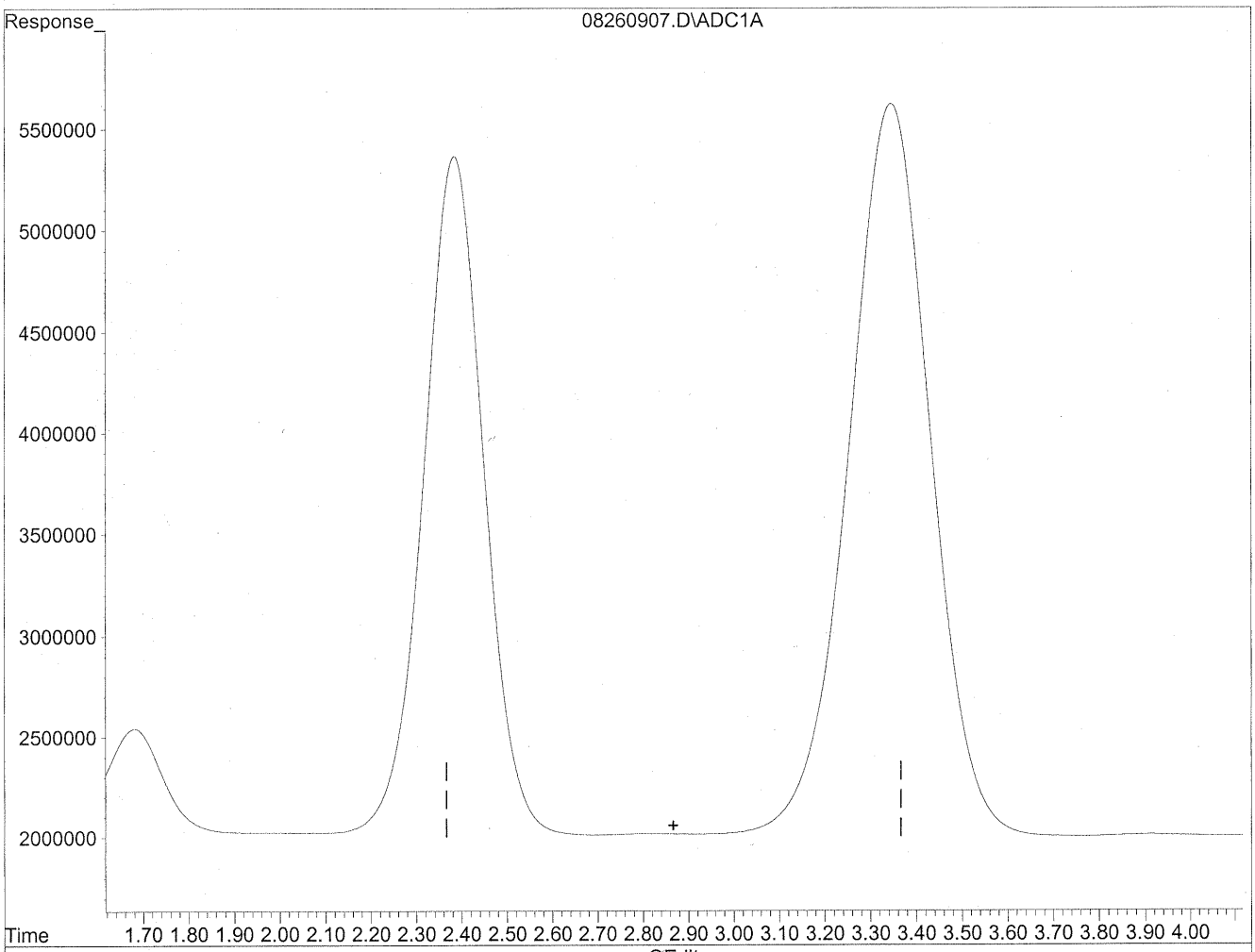
response 446740263

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260907.D Vial: 7
Acq On : 26 Aug 2009 6:35 pm Operator: HC
Sample : P0902942-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



QEdit

(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

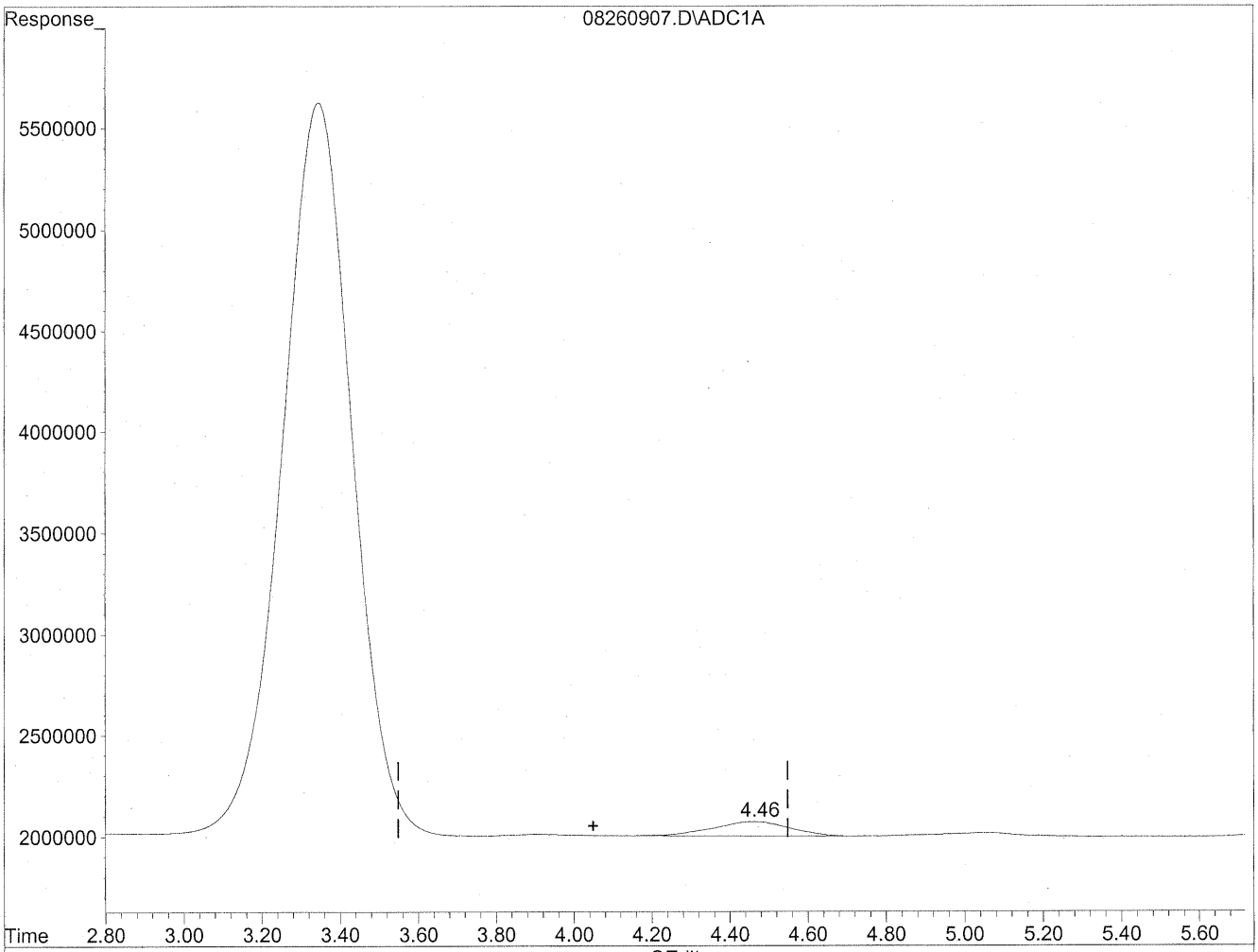
(+) = Expected Retention Time

08260907.D TO110709.M Sat Aug 29 16:47:44 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260907.D Vial: 7
Acq On : 26 Aug 2009 6:35 pm Operator: HC
Sample : P0902942-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.46min 97.924ng/ml

response 9539245

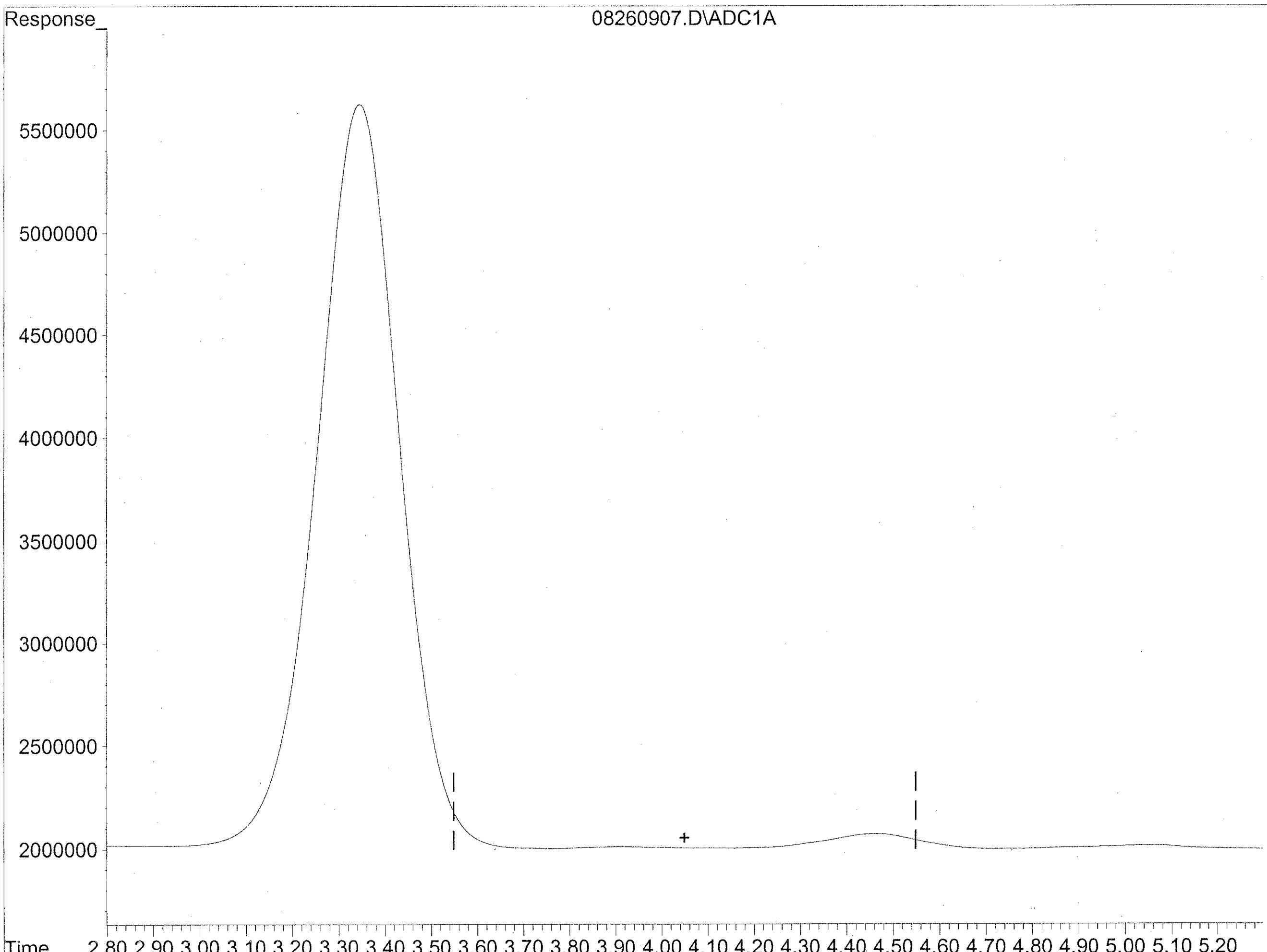
(+) = Expected Retention Time

08260907.D TO110709.M Sat Aug 29 16:47:47 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260907.D Vial: 7
Acq On : 26 Aug 2009 6:35 pm Operator: HC
Sample : P0902942-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



QEdit

(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/30/09 MP

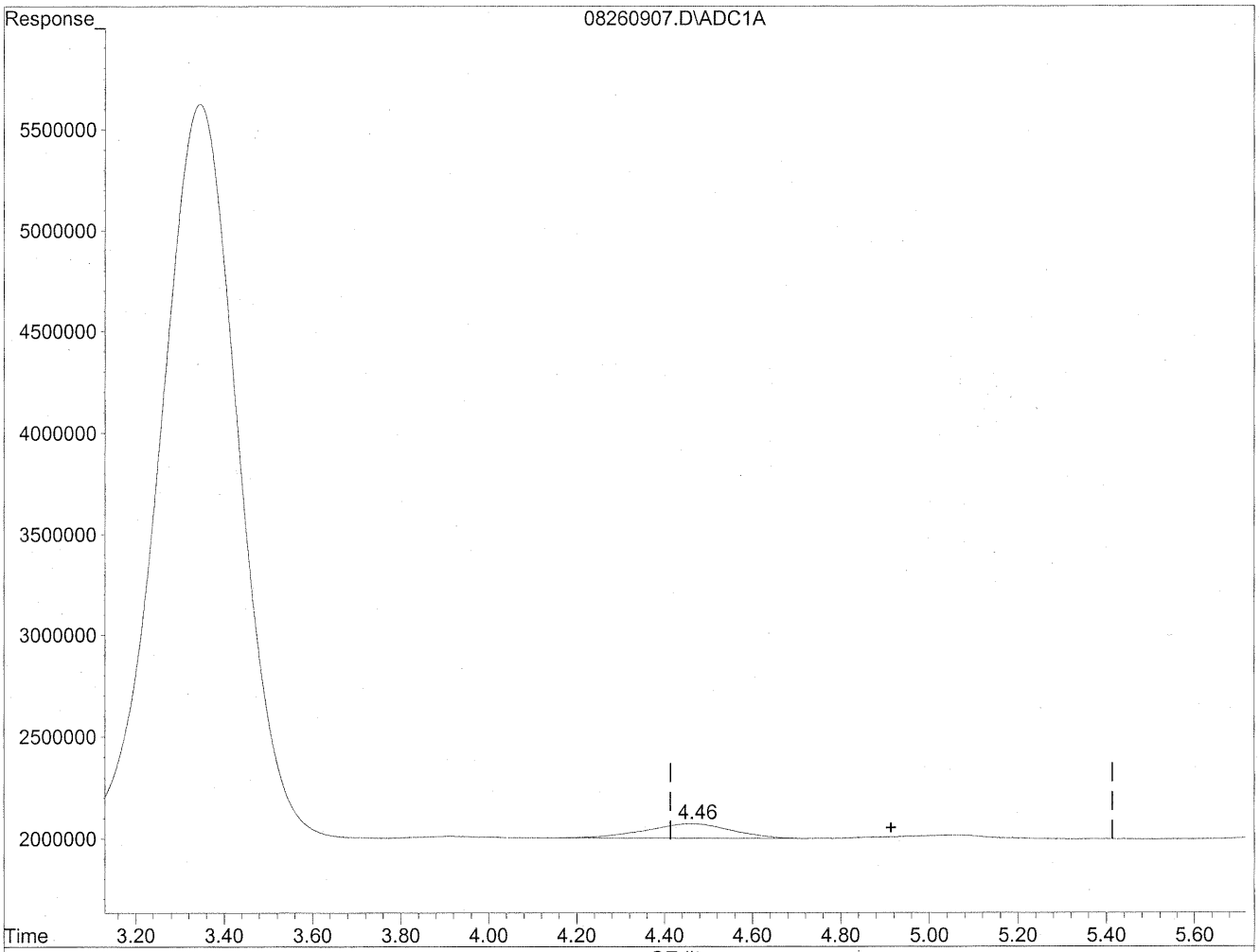
LW 8/31/09

(+) = Expected Retention Time
08260907.D TO110709.M Sat Aug 29 16:47:50 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260907.D Vial: 7
Acq On : 26 Aug 2009 6:35 pm Operator: HC
Sample : P0902942-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.46min 107.988ng/ml

response 9539245

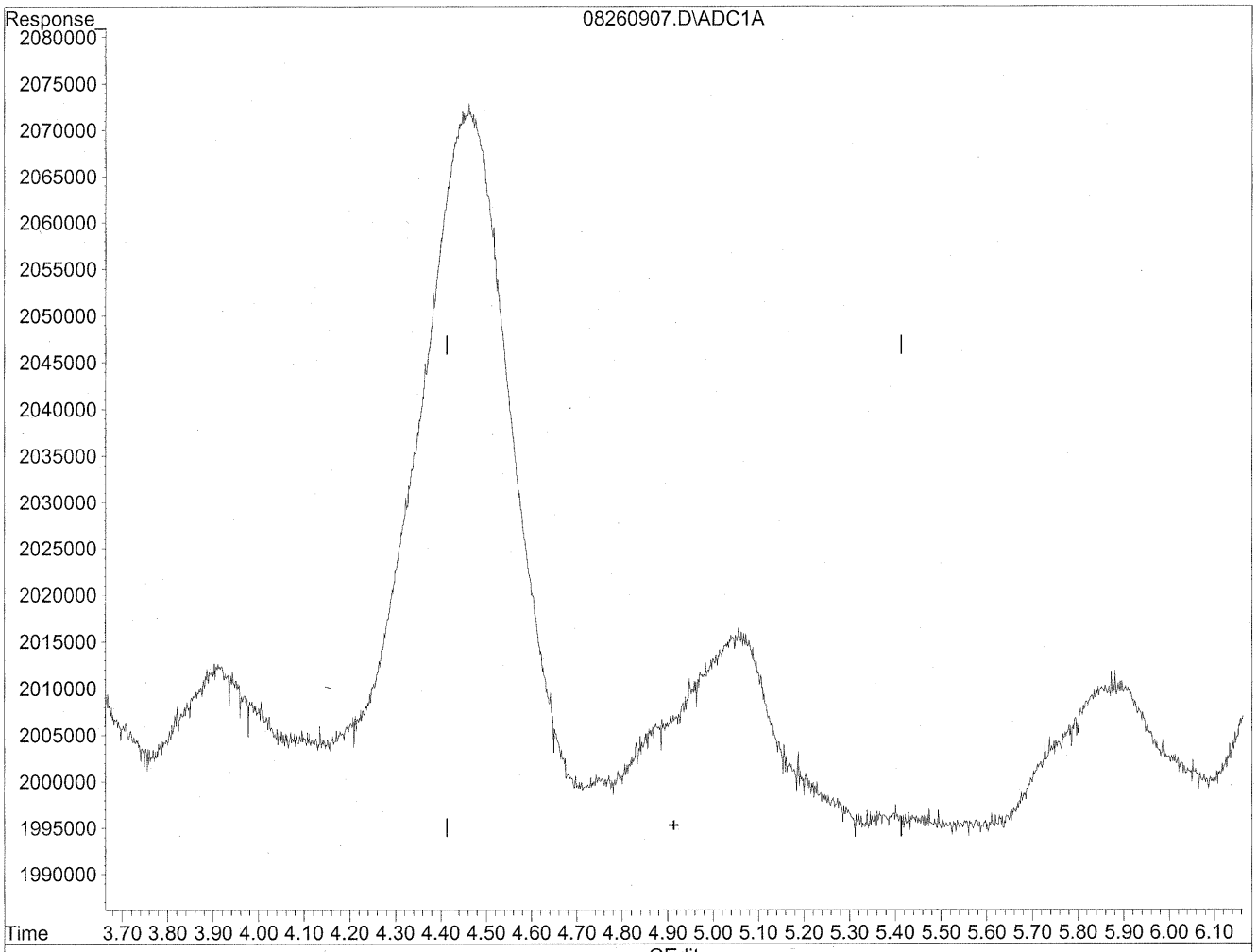
(+) = Expected Retention Time

08260907.D TO110709.M Sat Aug 29 16:47:55 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260907.D Vial: 7
Acq On : 26 Aug 2009 6:35 pm Operator: HC
Sample : P0902942-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

0.00min 0.000ng/ml d

response 0

HC 8/27/2009 MP

(+) = Expected Retention Time

08260907.D TO110709.M Sat Aug 29 16:47:58 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **101471**
Client Project ID: **16512**

CAS Project ID: P0902942
CAS Sample ID: P0902942-003

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: 8/24/09
Date Received: 8/25/09
Date Analyzed: 8/26-27/09
Desorption Volume: 1.0 ml
Volume Sampled: 100.5 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**200**	**2.0**	1.0	**1.6**	0.81	**BH**
75-07-0	**Acetaldehyde**	**110**	**1.1**	1.0	**0.63**	0.55	
123-38-6	Propionaldehyde	< 100	ND	1.0	ND	0.42	
4170-30-3	Crotonaldehyde, Total	< 200	ND	2.0	ND	0.69	**V**
123-72-8	Butyraldehyde	< 100	ND	1.0	ND	0.34	
100-52-7	Benzaldehyde	< 100	ND	1.0	ND	0.23	
590-86-3	Isovaleraldehyde	< 100	ND	1.0	ND	0.28	
110-62-3	Valeraldehyde	< 100	ND	1.0	ND	0.28	
529-20-4	o-Tolualdehyde	< 100	ND	1.0	ND	0.20	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	2.0	ND	0.41	
66-25-1	n-Hexaldehyde	< 100	ND	1.0	ND	0.24	
5779-94-2	2,5-Dimethylbenzaldehyde	< 200	ND	2.0	ND	0.36	**V**

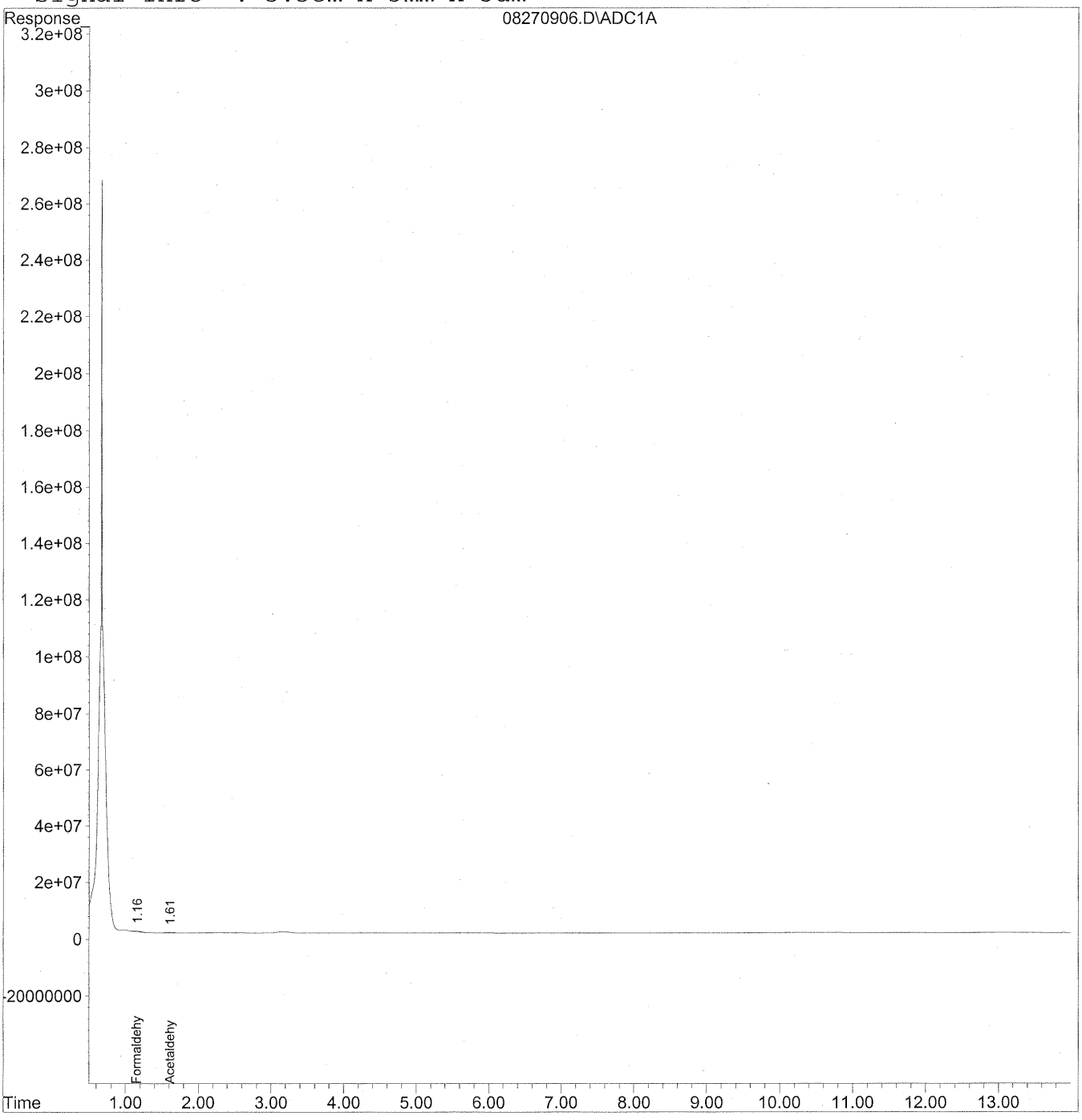
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
BH = The back section of the tube yielded higher results than the front.
V = The continuing calibration verification standard was outside (biased low) the specified limits for this compound.

Verified By: Re Date: 9/1/09

Data File : J:\LC01\DATA\TO11\2009_08\27\08270906.D Vial: 6
Acq On : 27 Aug 2009 10:20 am Operator: HC
Sample : P0902942-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 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\27\08270906.D Vial: 6
Acq On : 27 Aug 2009 10:20 am Operator: HC
Sample : P0902942-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 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

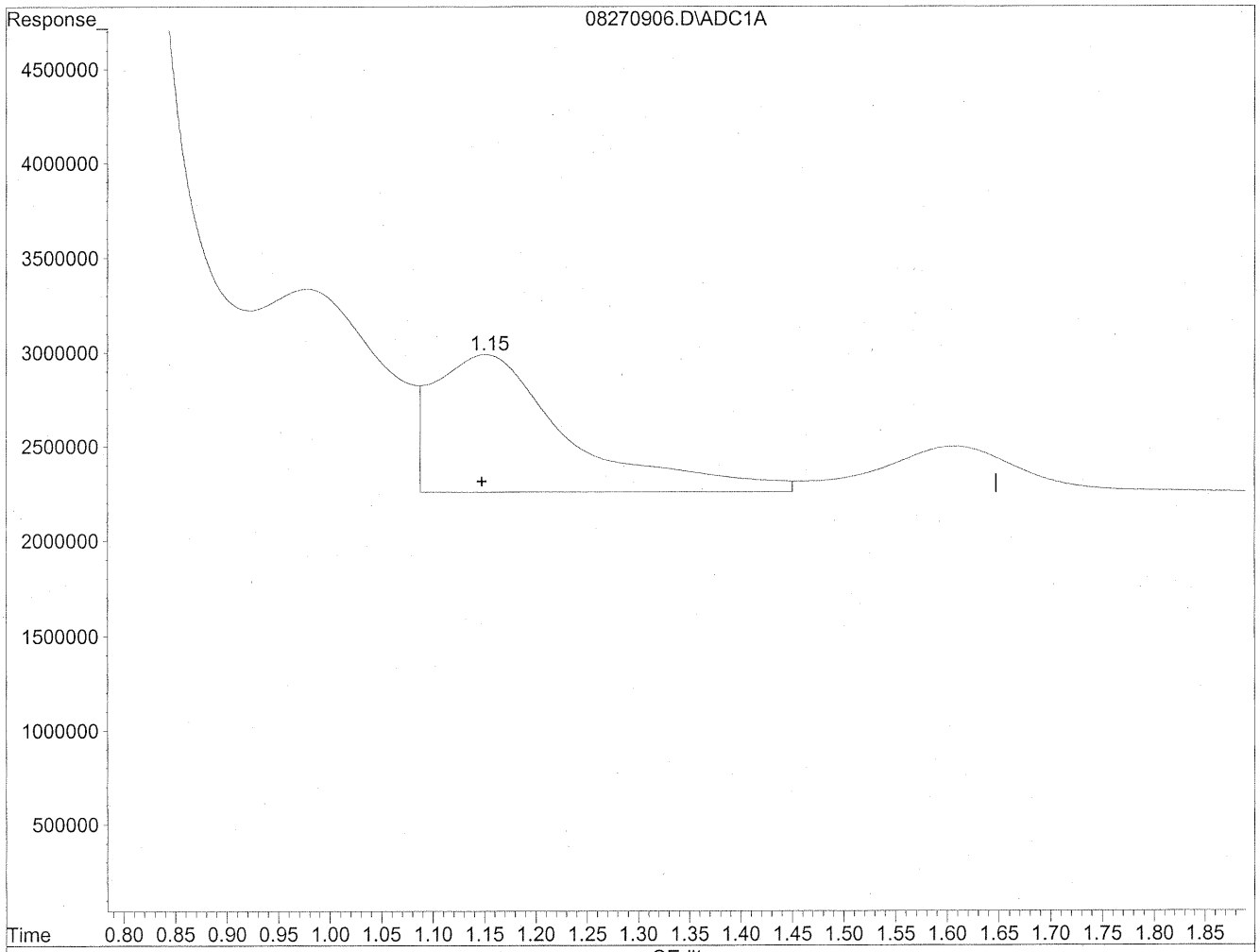
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	15666852	85.340	ng/mlm
2)	Acetaldehyde	1.61	15877696	113.231	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	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\27\08270906.D Vial: 6
Acq On : 27 Aug 2009 10:20 am Operator: HC
Sample : P0902942-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:32 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.15min 363.903ng/ml

response 66805820

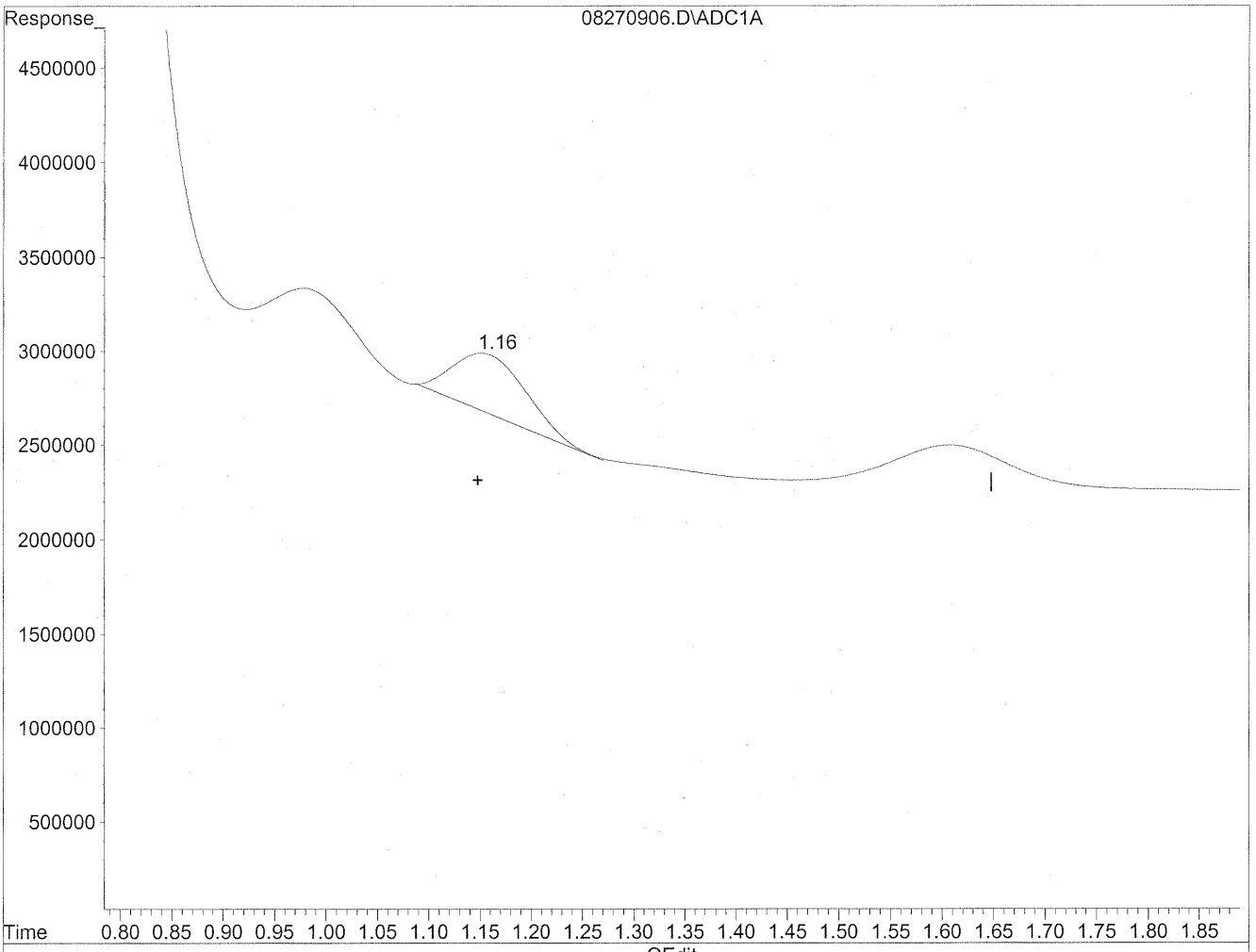
(+) = Expected Retention Time

08270906.D TO110709.M Sun Aug 30 13:32:30 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270906.D Vial: 6
Acq On : 27 Aug 2009 10:20 am Operator: HC
Sample : P0902942-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:32 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.16min 85.340ng/ml m

response 15666852

8/31/09

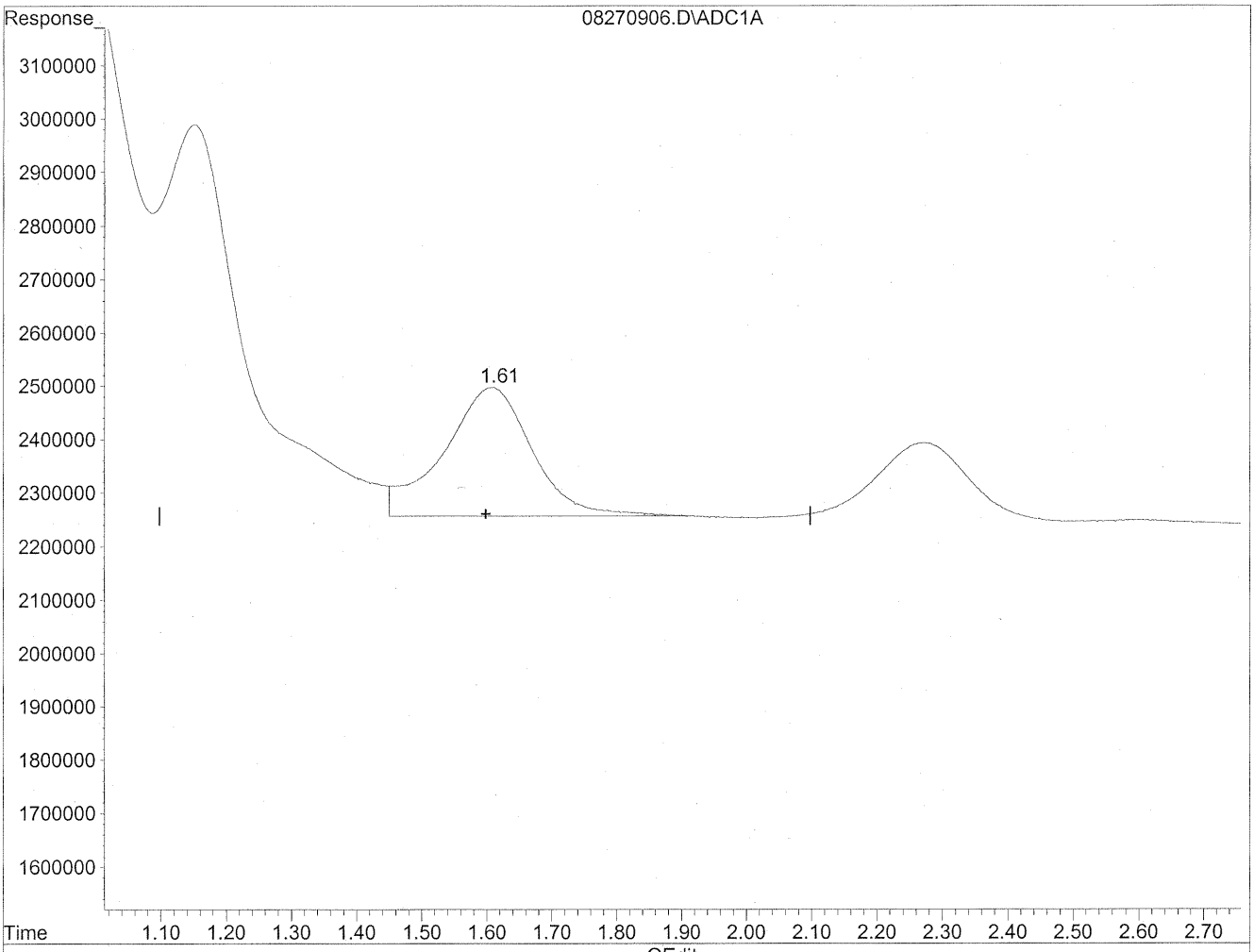
8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270906.D Vial: 6
Acq On : 27 Aug 2009 10:20 am Operator: HC
Sample : P0902942-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:32 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.61min 162.254ng/ml

response 22751796

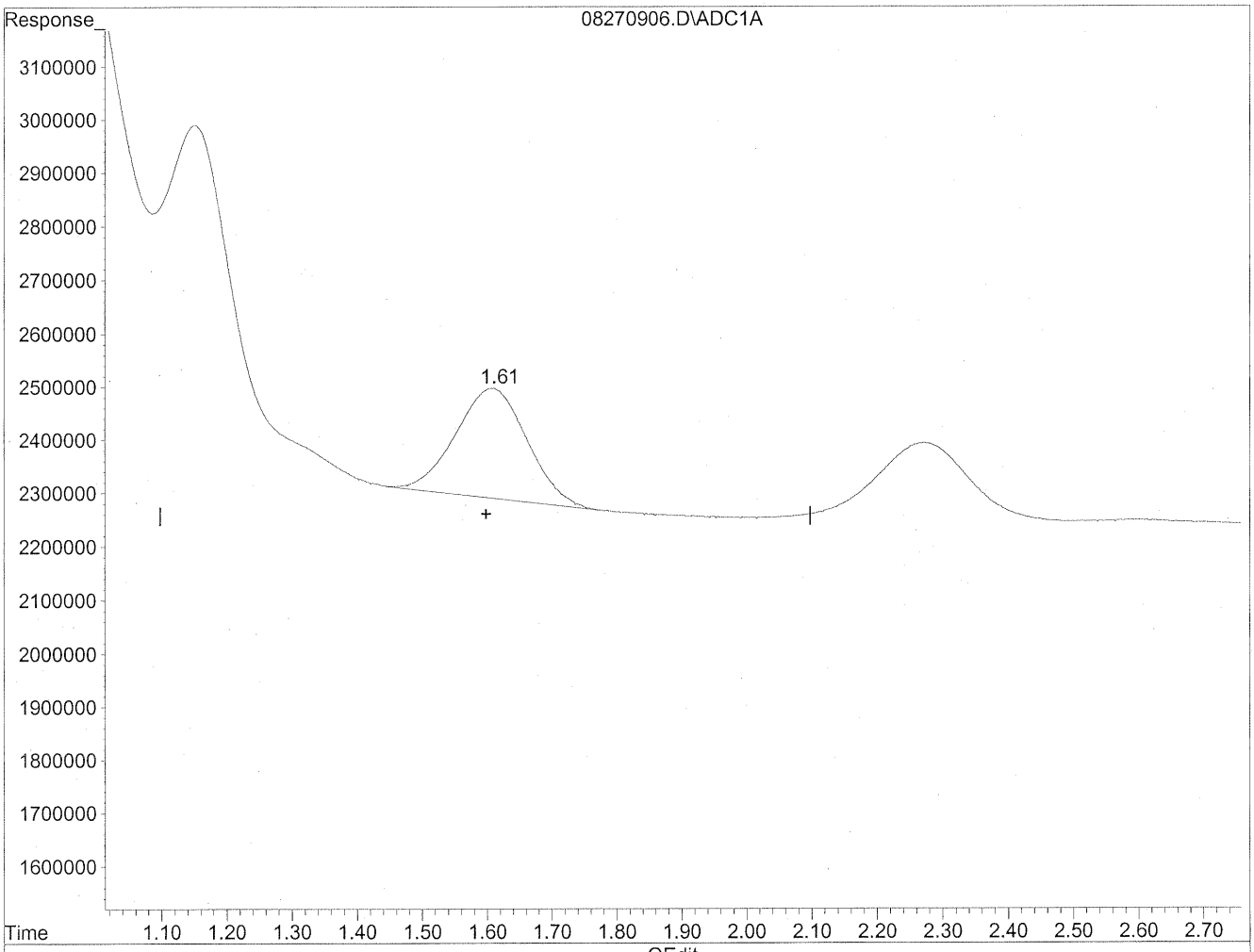
(+) = Expected Retention Time

08270906.D TO110709.M Sun Aug 30 13:32:42 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270906.D Vial: 6
Acq On : 27 Aug 2009 10:20 am Operator: HC
Sample : P0902942-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:32 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.61min 113.231ng/ml m

response 15877696

HC 8/31/09 1C

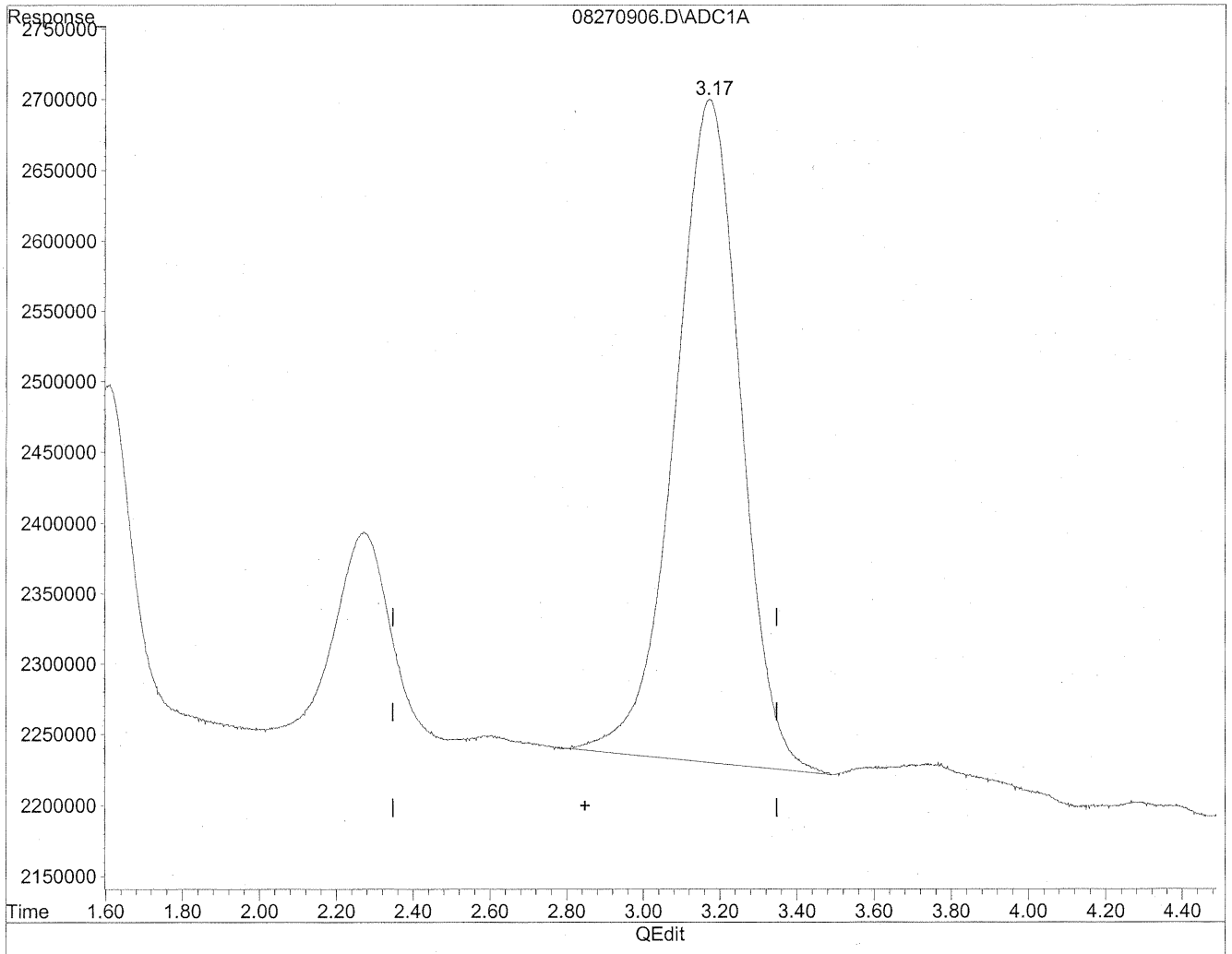
LW 8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270906.D Vial: 6
Acq On : 27 Aug 2009 10:20 am Operator: HC
Sample : P0902942-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:32 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.17min 527.335ng/ml

response 56264118

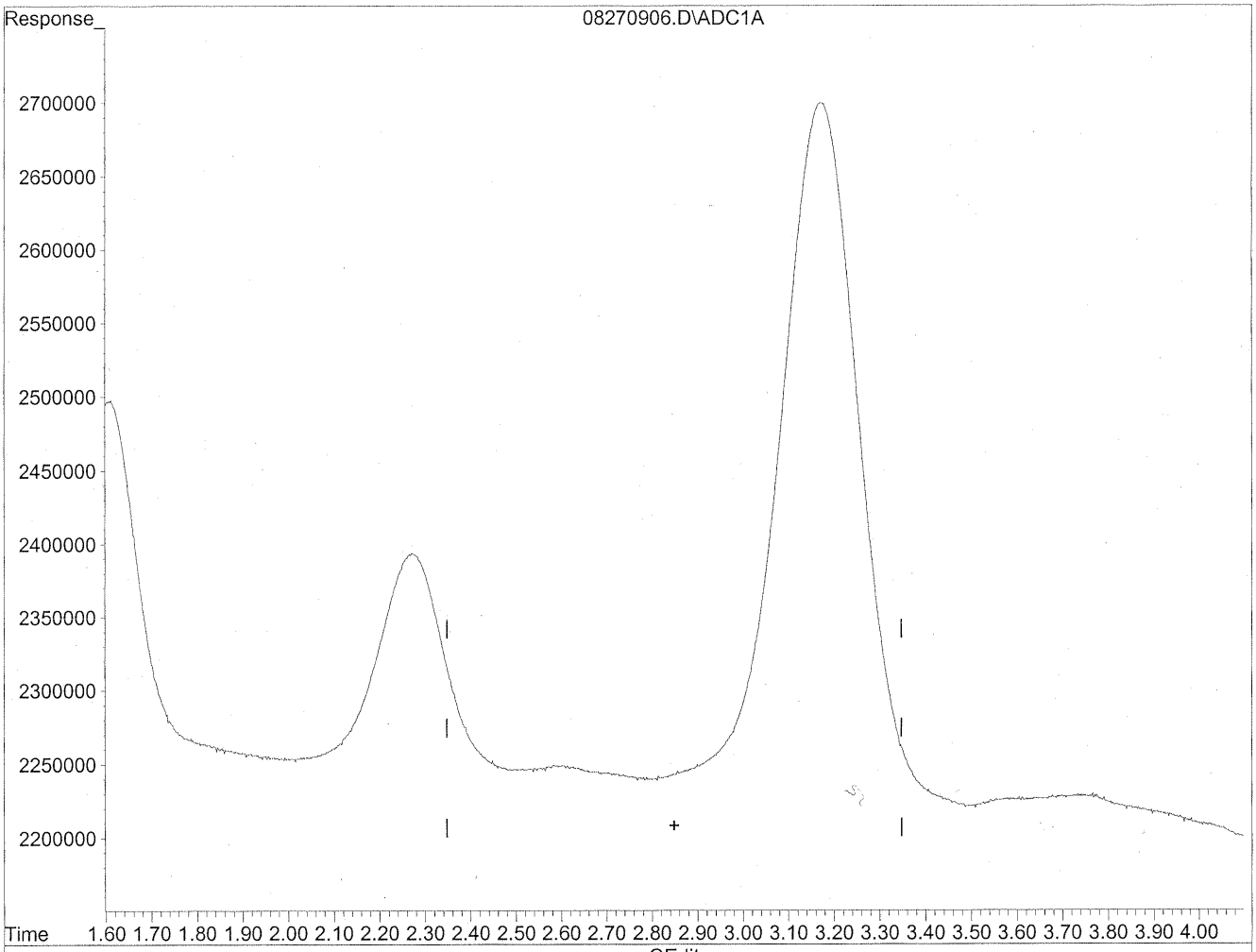
(+) = Expected Retention Time

08270906.D TO110709.M Sun Aug 30 13:32:59 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270906.D Vial: 6
Acq On : 27 Aug 2009 10:20 am Operator: HC
Sample : P0902942-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:32 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/31/09 WP

8/31/09

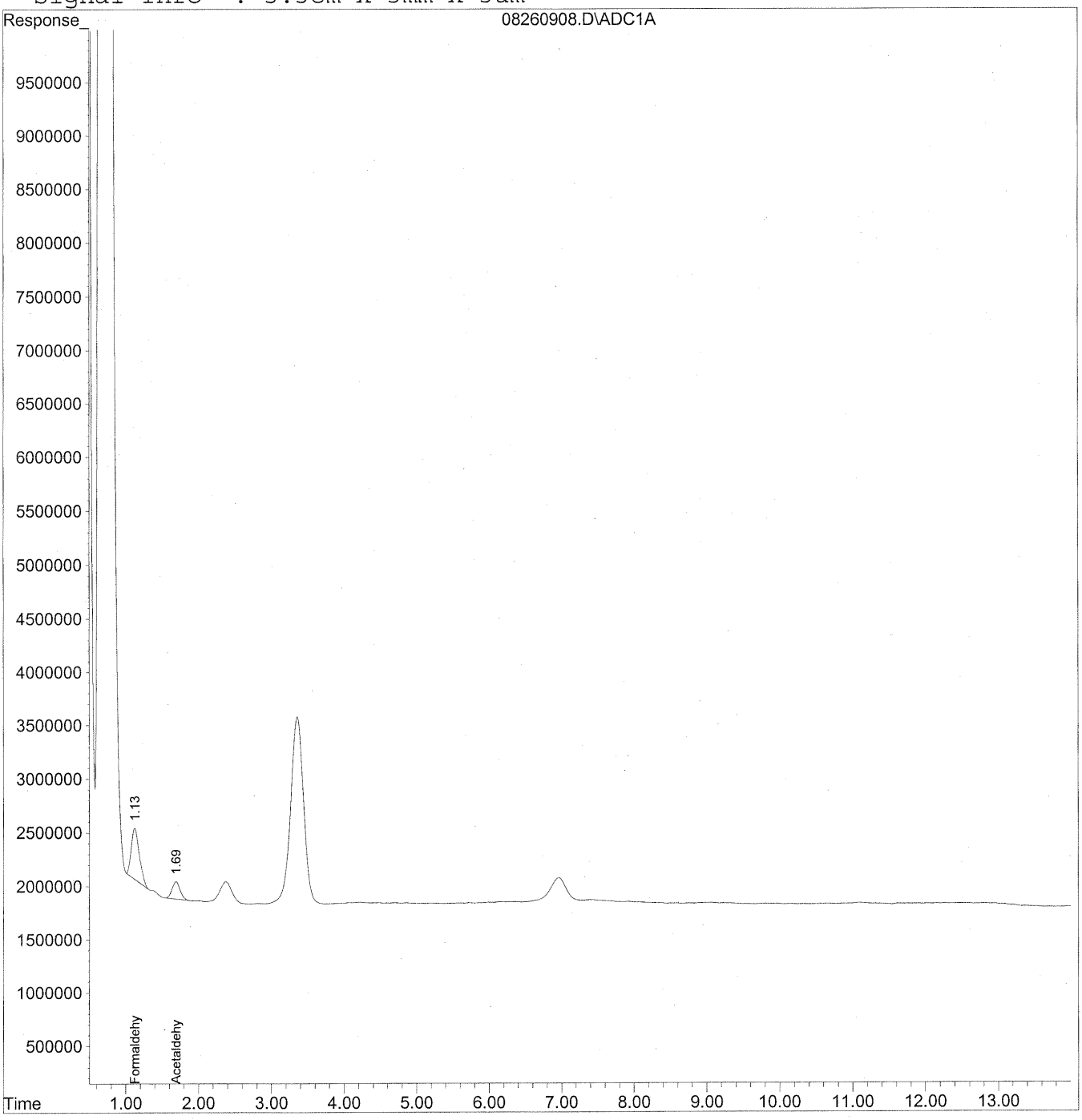
75

(+) = Expected Retention Time
08270906.D TO110709.M Sun Aug 30 13:33:02 2009

Data File : J:\LC01\DATA\TO11\2009_08\26\08260908.D Vial: 8
Acq On : 26 Aug 2009 6:50 pm Operator: HC
Sample : P0902942-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:48 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 : Thu Aug 27 07:35:56 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\26\08260908.D Vial: 8
Acq On : 26 Aug 2009 6:50 pm Operator: HC
Sample : P0902942-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:48 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 : Thu Aug 27 07:35:56 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

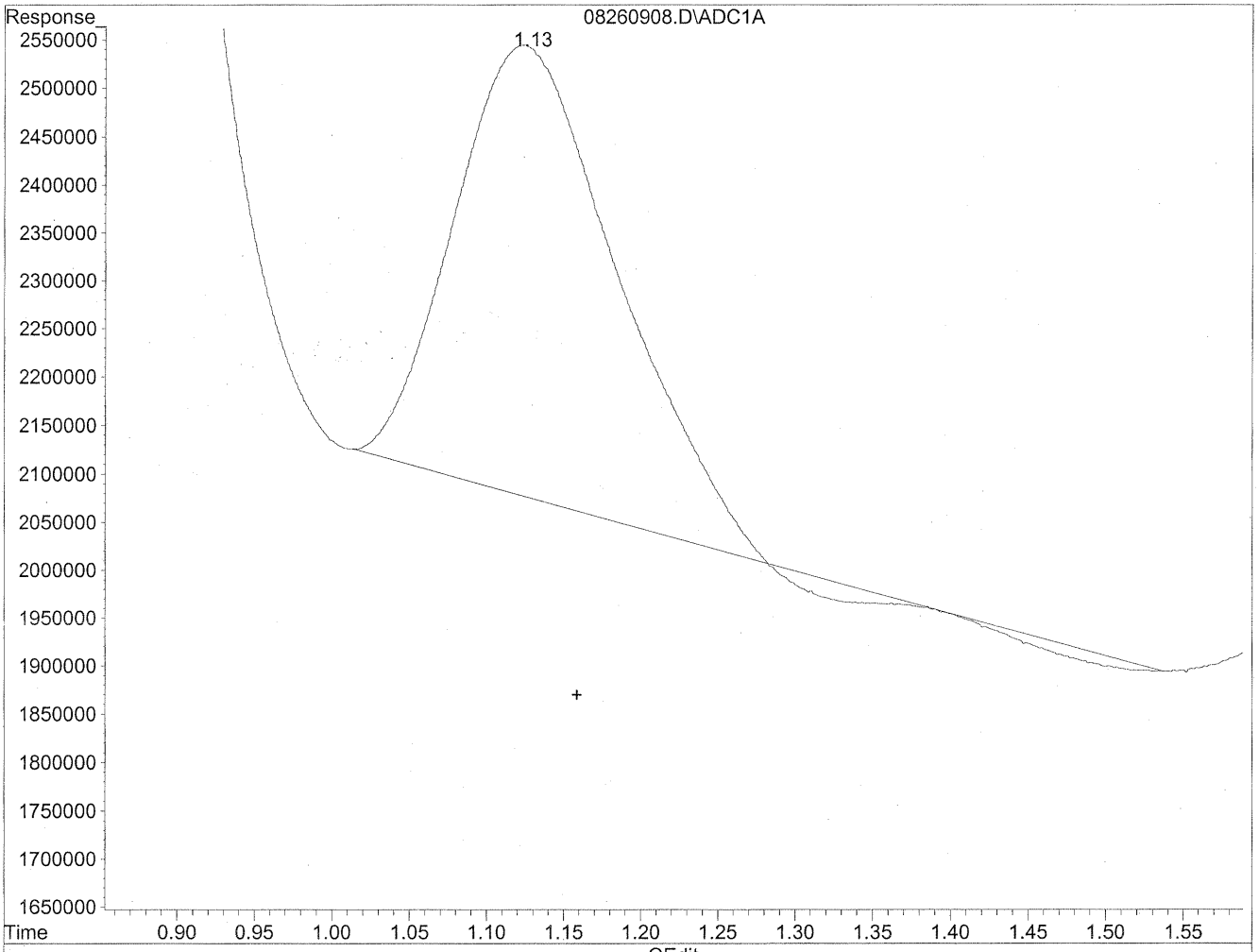
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.13	36474536	198.683	ng/mlm
2)	Acetaldehyde	1.69	12812934	91.375	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/mld
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\26\08260908.D Vial: 8
Acq On : 26 Aug 2009 6:50 pm Operator: HC
Sample : P0902942-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.12min 182.879ng/ml

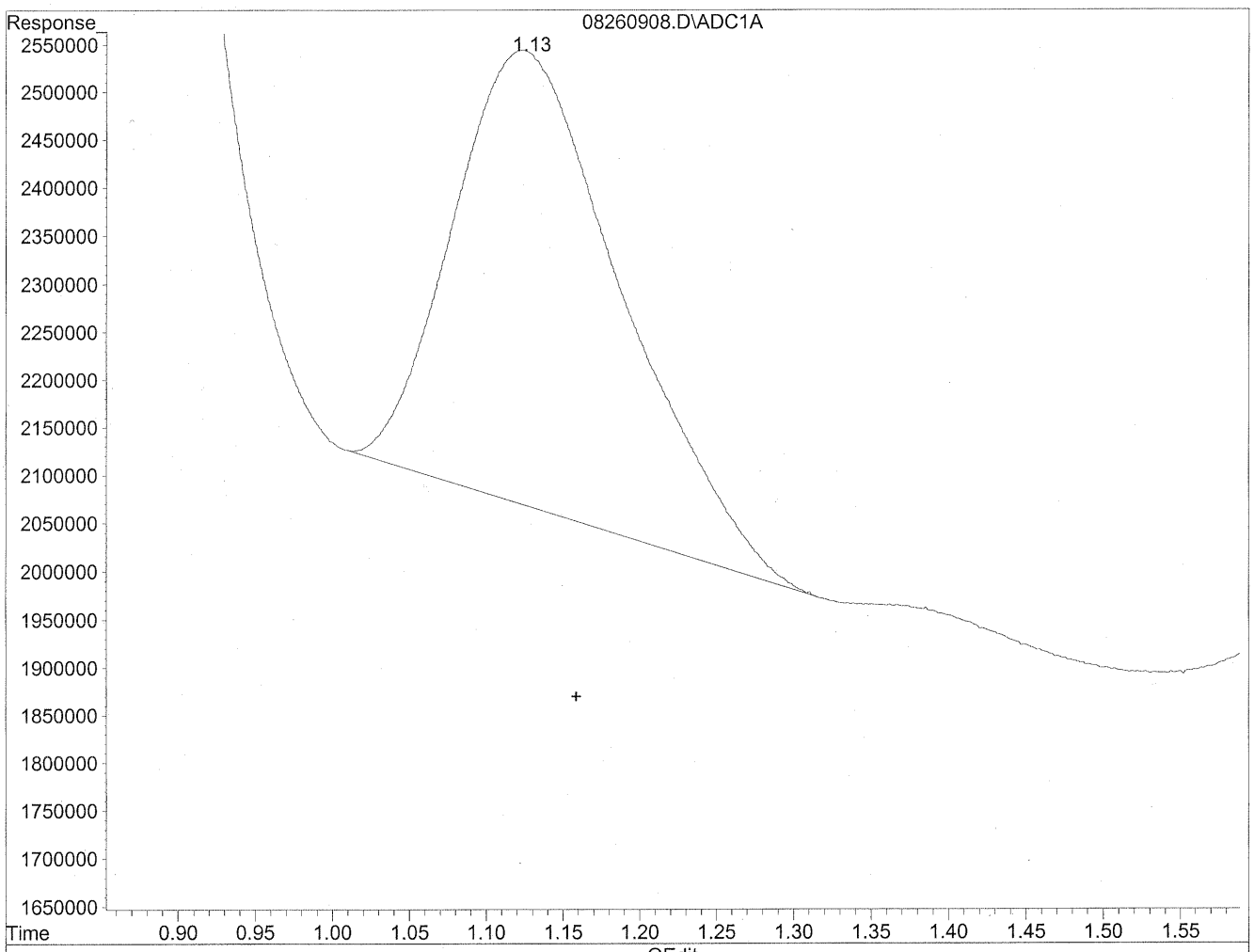
response 33573206

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260908.D Vial: 8
Acq On : 26 Aug 2009 6:50 pm Operator: HC
Sample : P0902942-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.13min 198.683ng/ml m

response 36474536

HC 8/30/09 IC

WP 9/3/09

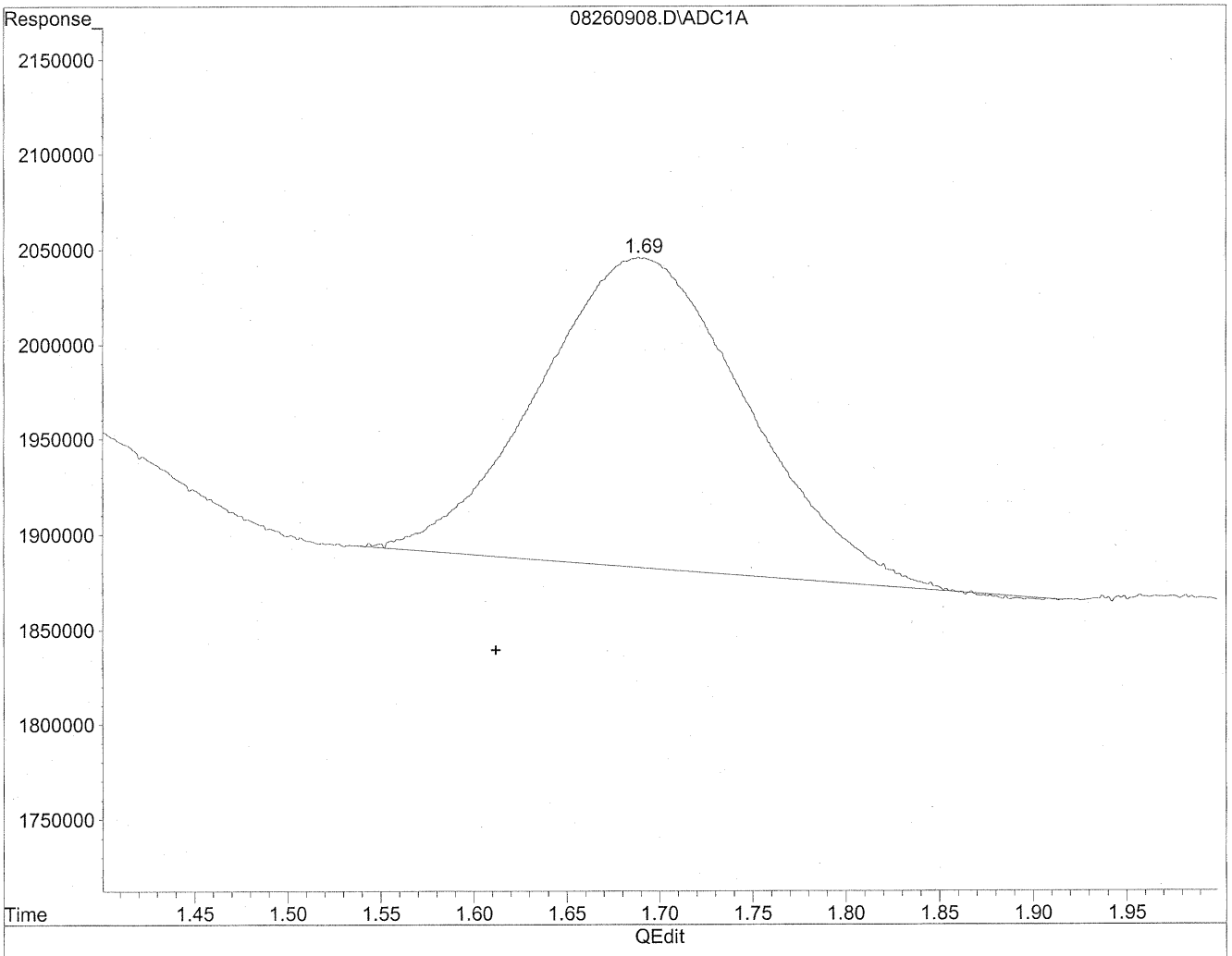
79

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260908.D Vial: 8
Acq On : 26 Aug 2009 6:50 pm Operator: HC
Sample : P0902942-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.69min 90.544ng/ml

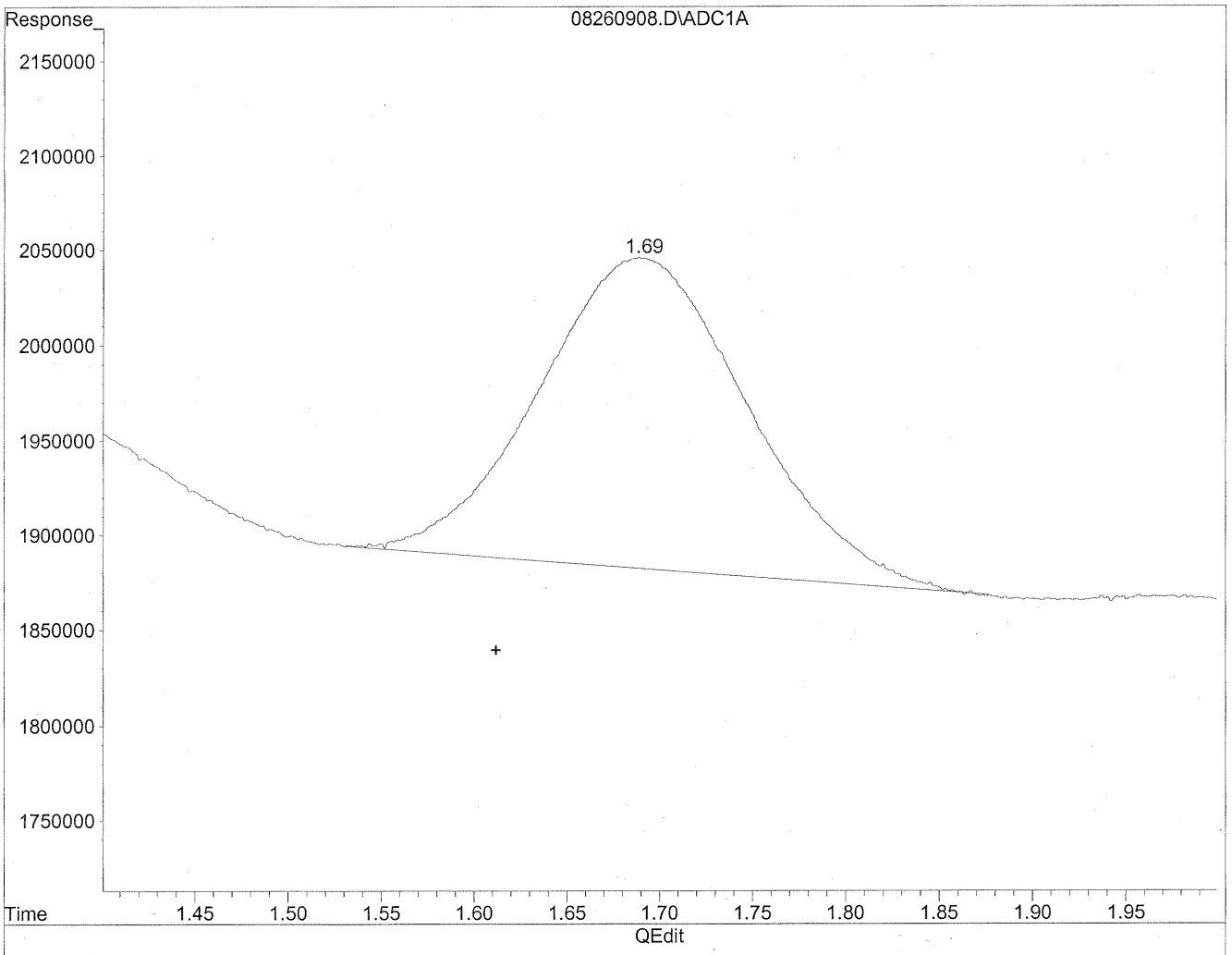
response 12696354

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260908.D Vial: 8
Acq On : 26 Aug 2009 6:50 pm Operator: HC
Sample : P0902942-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.69min 91.375ng/ml m

response 12812934

HC 8/31/09

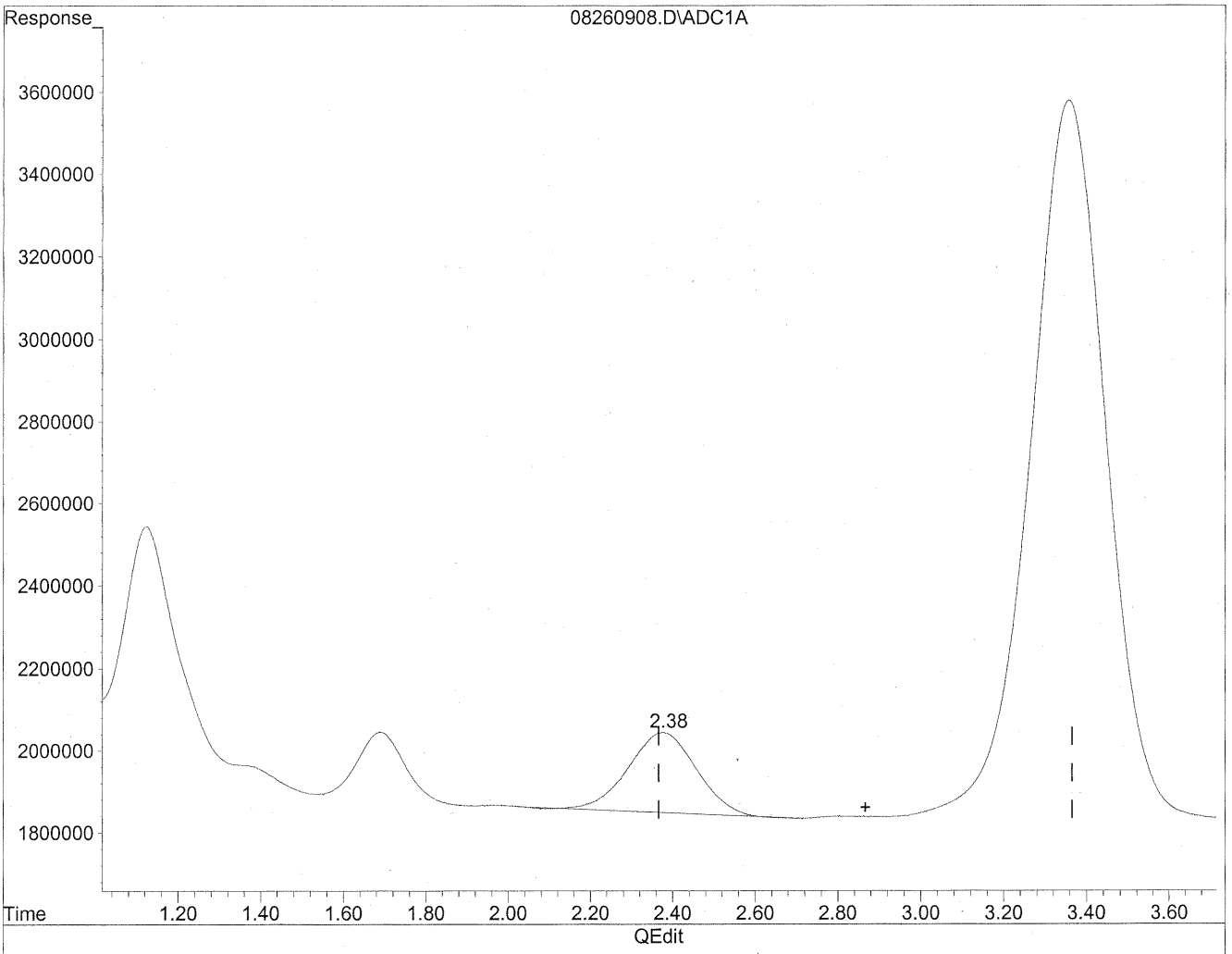
WK 8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260908.D Vial: 8
Acq On : 26 Aug 2009 6:50 pm Operator: HC
Sample : P0902942-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.38min 209.669ng/ml

response 22370651

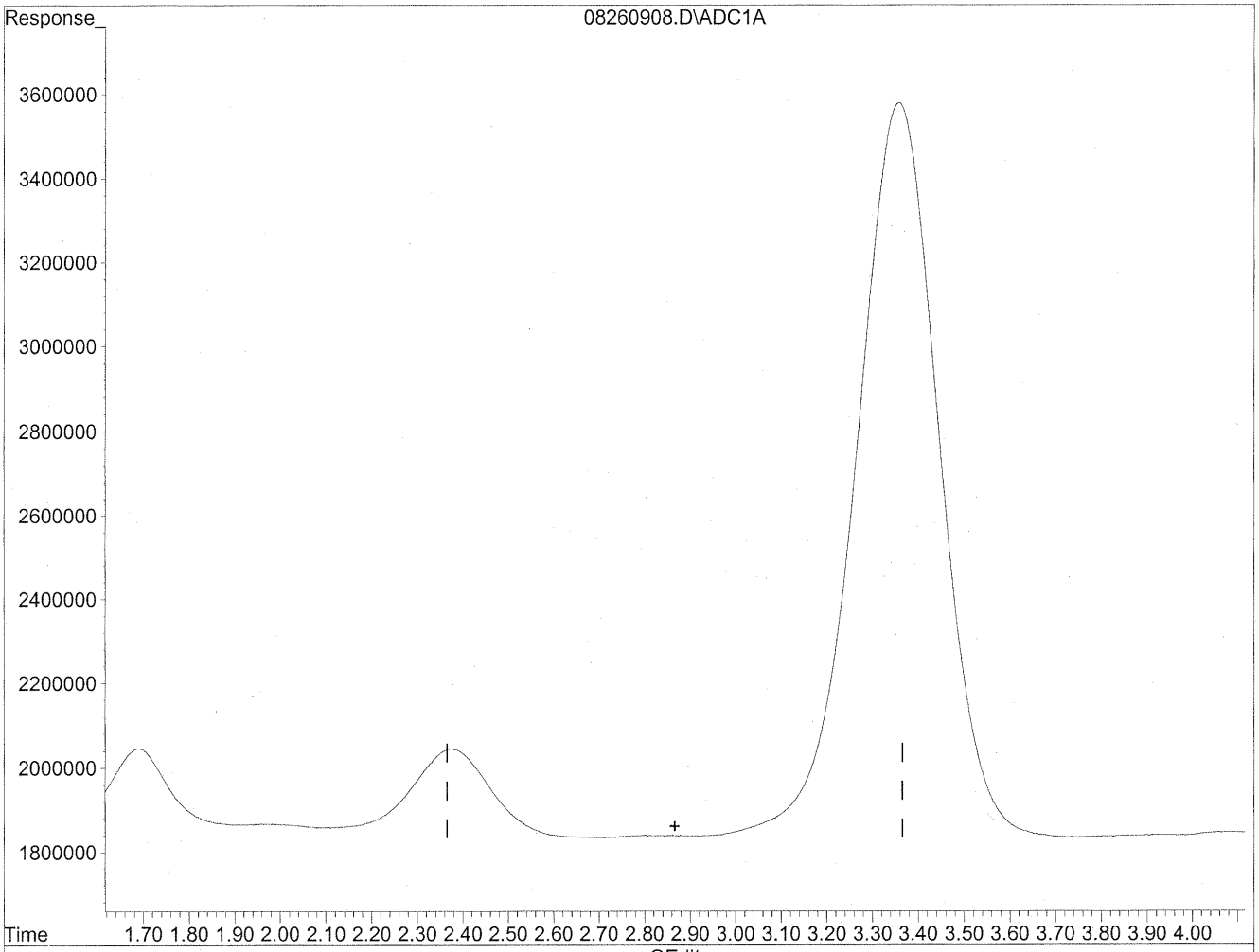
(+) = Expected Retention Time

08260908.D TO110709.M Sat Aug 29 16:48:43 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260908.D Vial: 8
Acq On : 26 Aug 2009 6:50 pm Operator: HC
Sample : P0902942-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/30/09 mP

WH 8/31/09

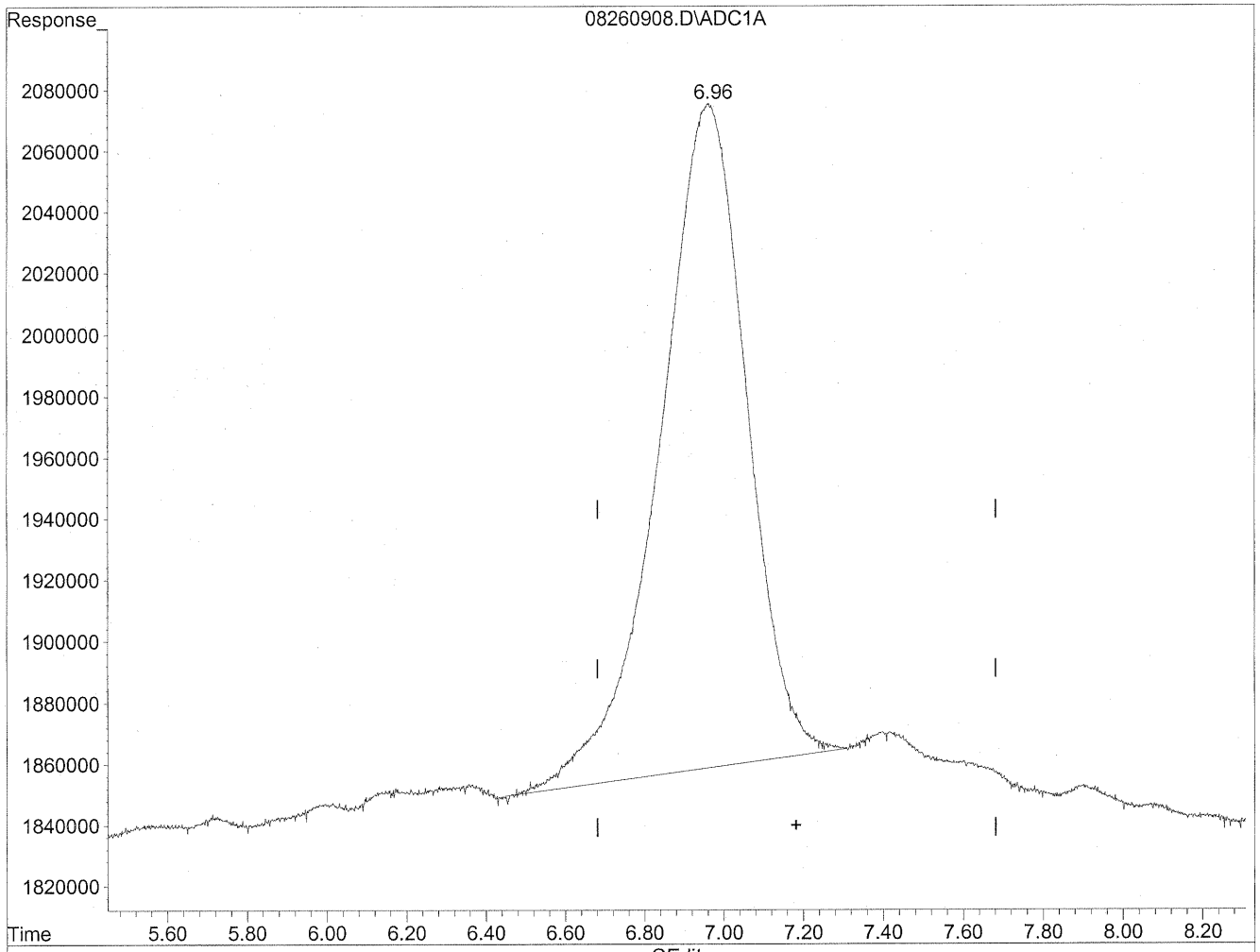
(+) = Expected Retention Time

08260908.D TO110709.M Sat Aug 29 16:48:49 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260908.D Vial: 8
Acq On : 26 Aug 2009 6:50 pm Operator: HC
Sample : P0902942-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

6.96min 421.415ng/ml

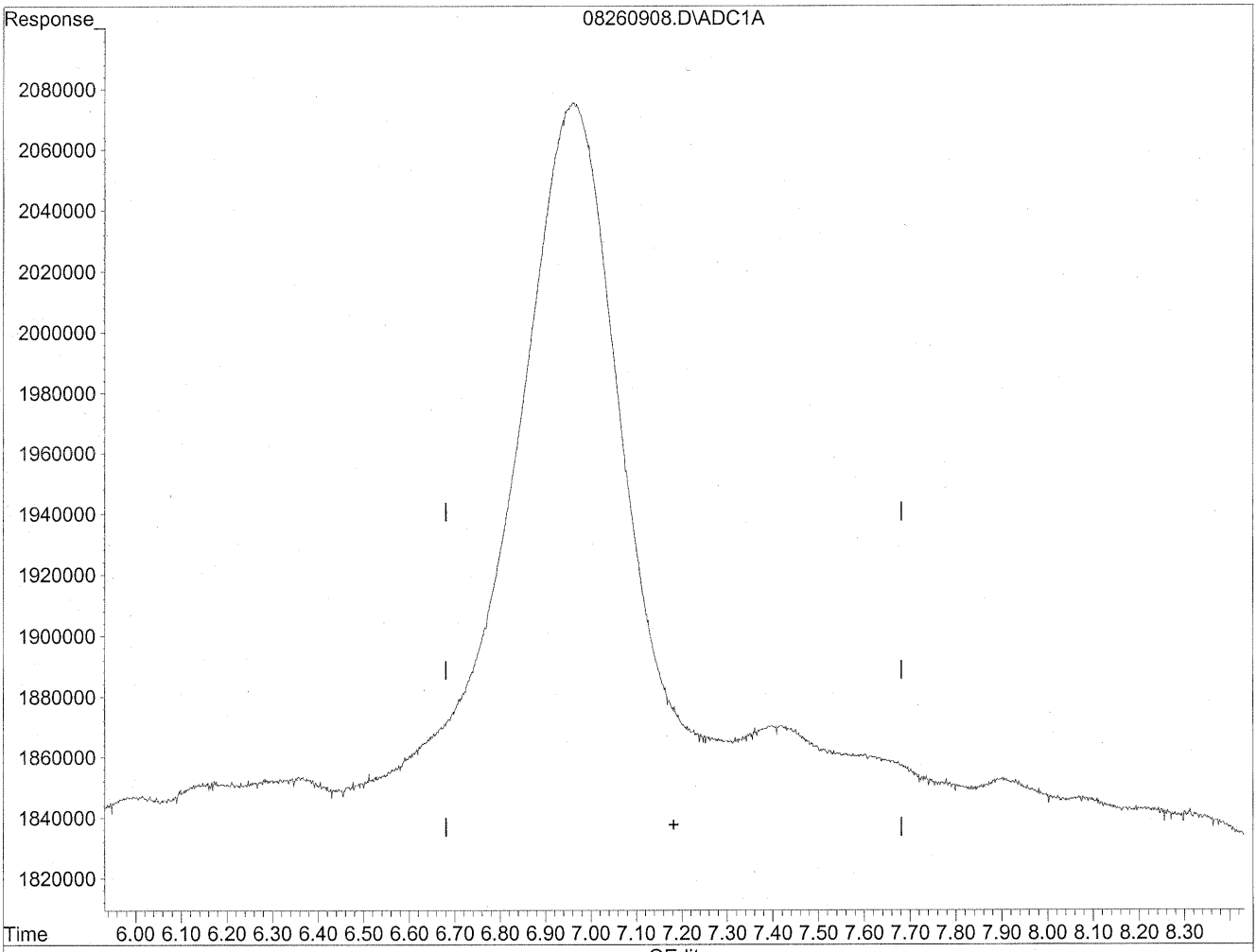
response 32976114

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260908.D Vial: 8
Acq On : 26 Aug 2009 6:50 pm Operator: HC
Sample : P0902942-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

0.00min 0.000ng/ml d

response 0

HC 8/30/09 MP

8/31/09

(+) = Expected Retention Time

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **101472**
Client Project ID: **16512**

CAS Project ID: P0902942
CAS Sample ID: P0902942-004

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: 8/24/09
Date Received: 8/25/09
Date Analyzed: 8/26-27/09
Desorption Volume: 1.0 ml
Volume Sampled: 101.5 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**2,100**	**21**	0.99	**17**	0.80	
75-07-0	**Acetaldehyde**	**4,300**	**43**	0.99	**24**	0.55	
123-38-6	**Propionaldehyde**	**540**	**5.3**	0.99	**2.2**	0.41	
4170-30-3	Crotonaldehyde, Total	< 200	ND	2.0	ND	0.69	**V**
123-72-8	**Butyraldehyde**	**1,000**	**9.9**	0.99	**3.4**	0.33	
100-52-7	**Benzaldehyde**	**820**	**8.1**	0.99	**1.9**	0.23	
590-86-3	**Isovaleraldehyde**	**170**	**1.7**	0.99	**0.47**	0.28	
110-62-3	**Valeraldehyde**	**1,400**	**14**	0.99	**4.0**	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**	**4,200**	**41**	0.99	**10**	0.24	
5779-94-2	2,5-Dimethylbenzaldehyde	< 200	ND	2.0	ND	0.36	**V**

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.
V = The continuing calibration verification standard was outside (biased low) the specified limits for this compound.

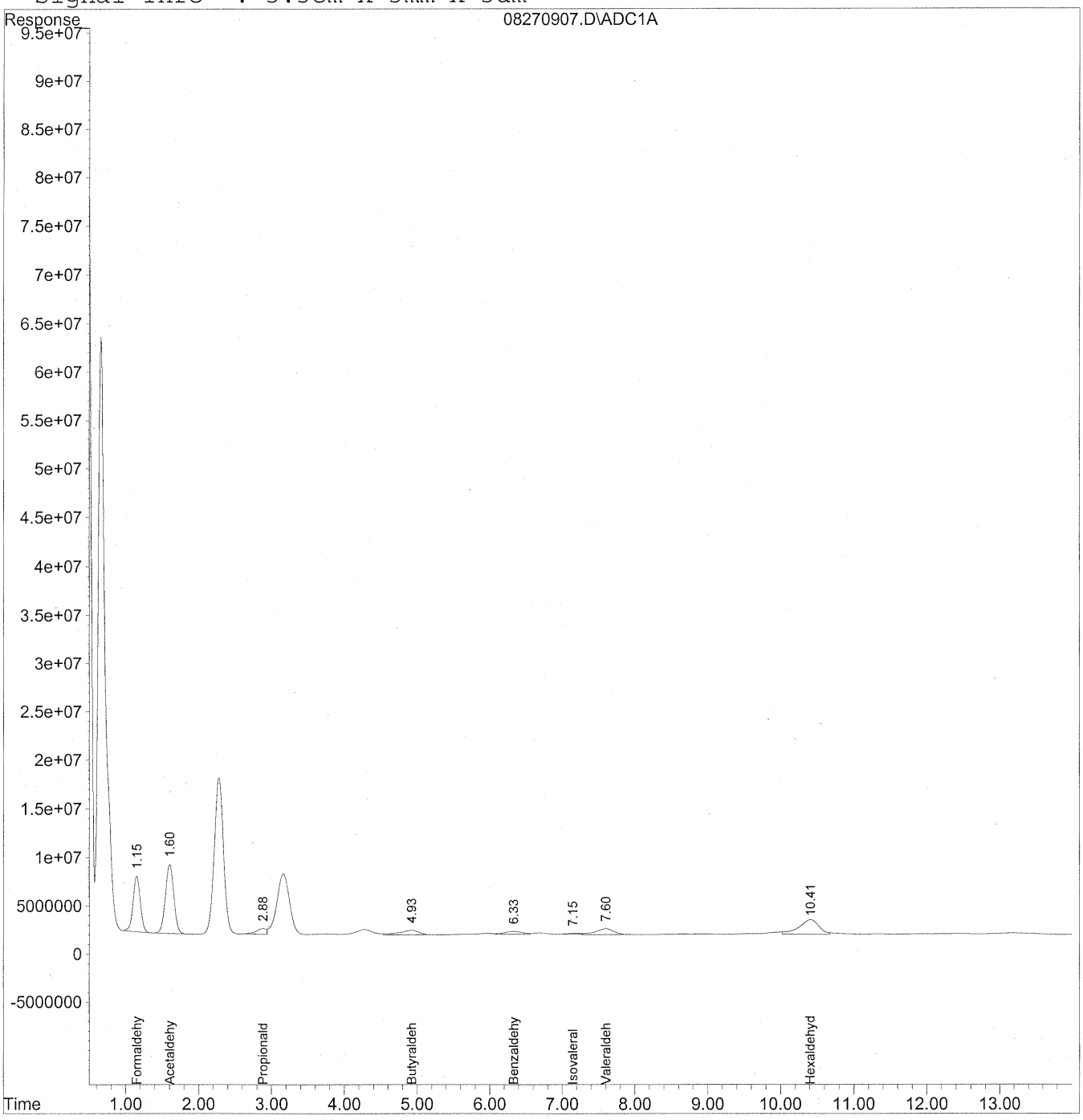
Verified By: RC Date: 9/11/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:36 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 : Sat Aug 29 17:49:00 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



87

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:36 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 : Sat Aug 29 17:49:00 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

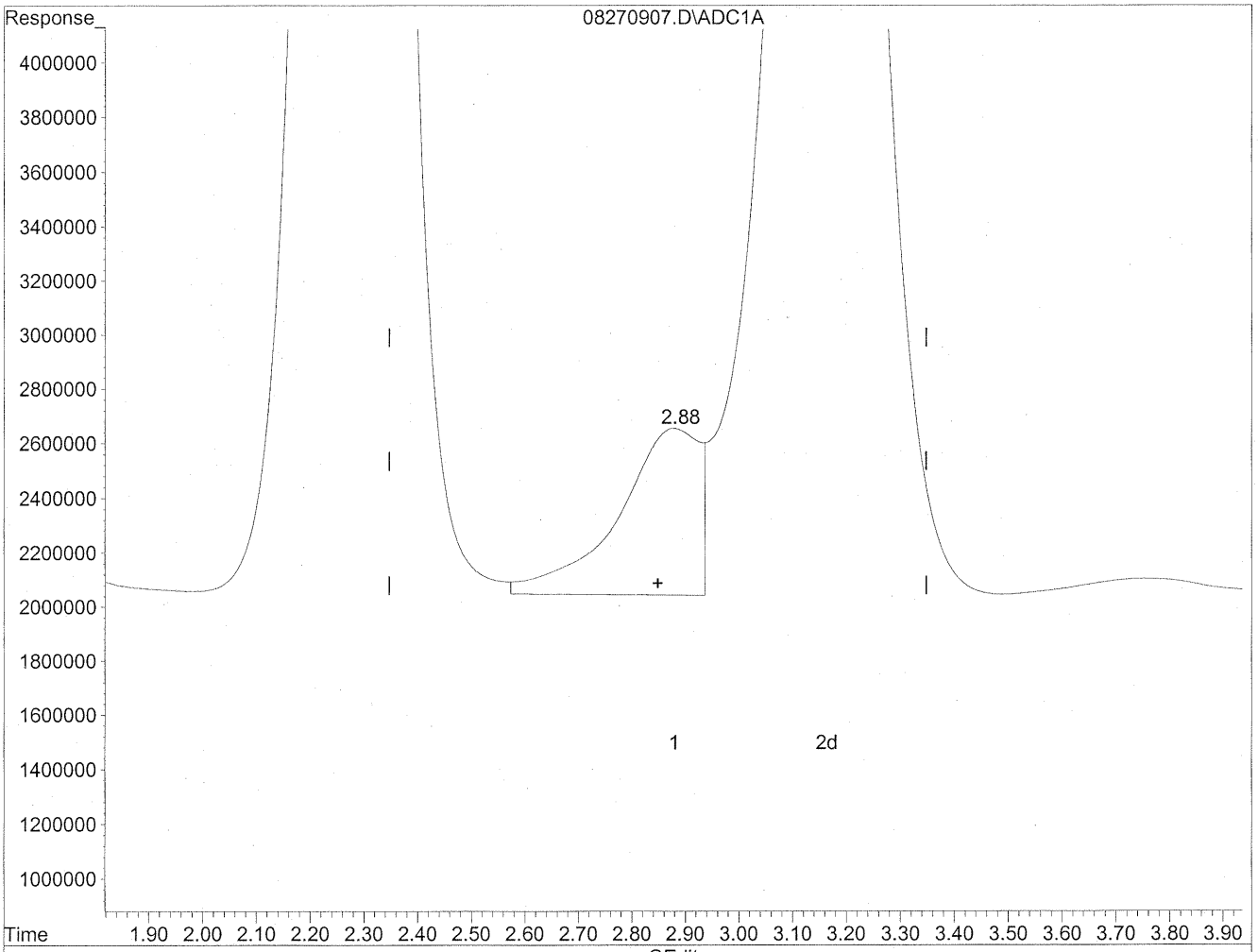
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.15	389092830	2119.457	ng/ml
2)	Acetaldehyde	1.60	571008109	4072.128	ng/ml
3)	Propionaldehyde	2.88	57601393	539.869	ng/mlm
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	4.93	89011049	1007.640	ng/mlm
6)	Benzaldehyde	6.33	54017560	820.072	ng/mlm
7)	Isovaleraldehyde	7.15	13119955	167.665	ng/mlm
8)	Valeraldehyde	7.60	105920278	1440.994	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	10.41	282428395	4193.832	ng/mlm
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/mld

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.88min 602.983ng/ml

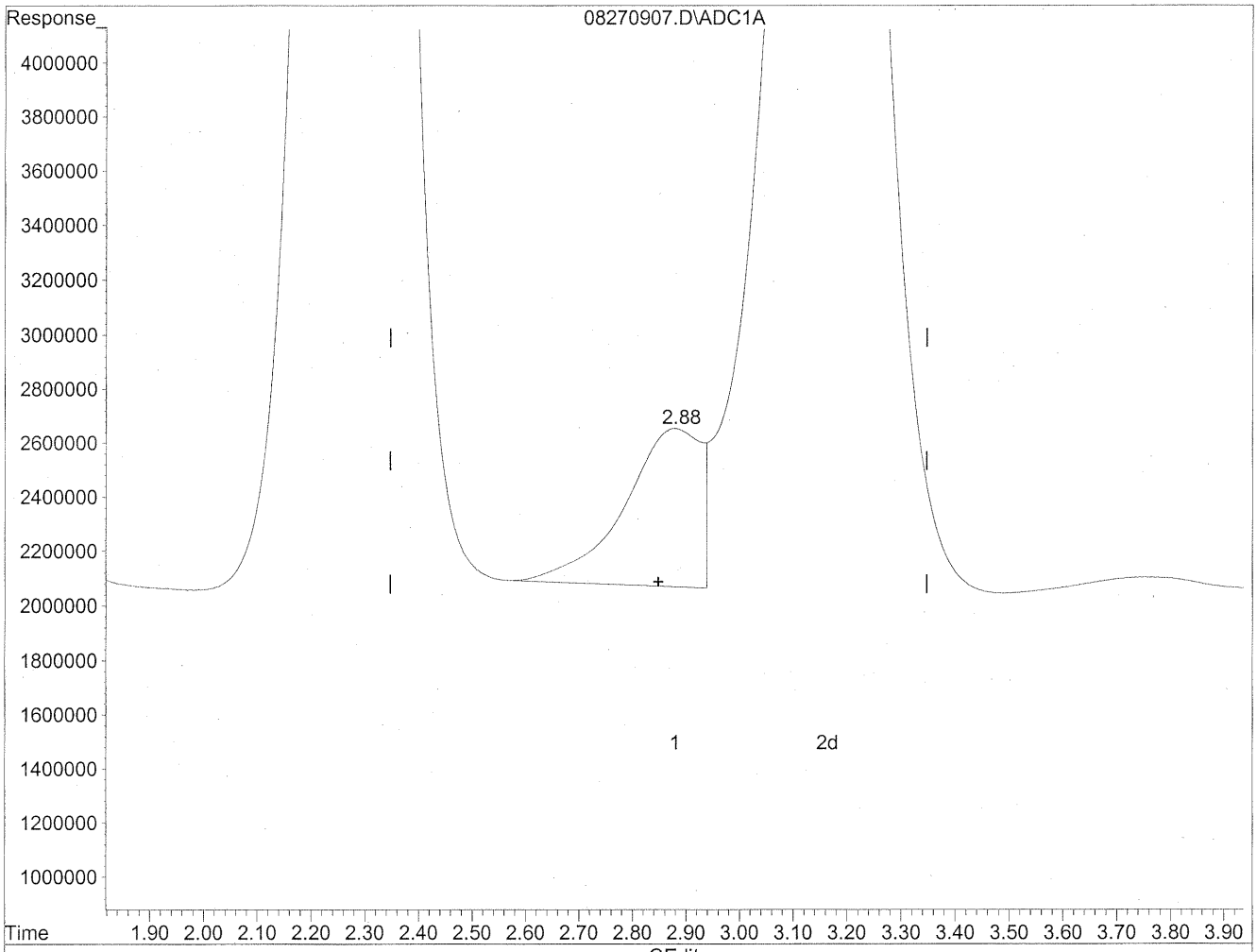
response 64335447

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.88min 539.869ng/ml m

response 57601393

JC 8/31/09

BC

WDS 8/31/09

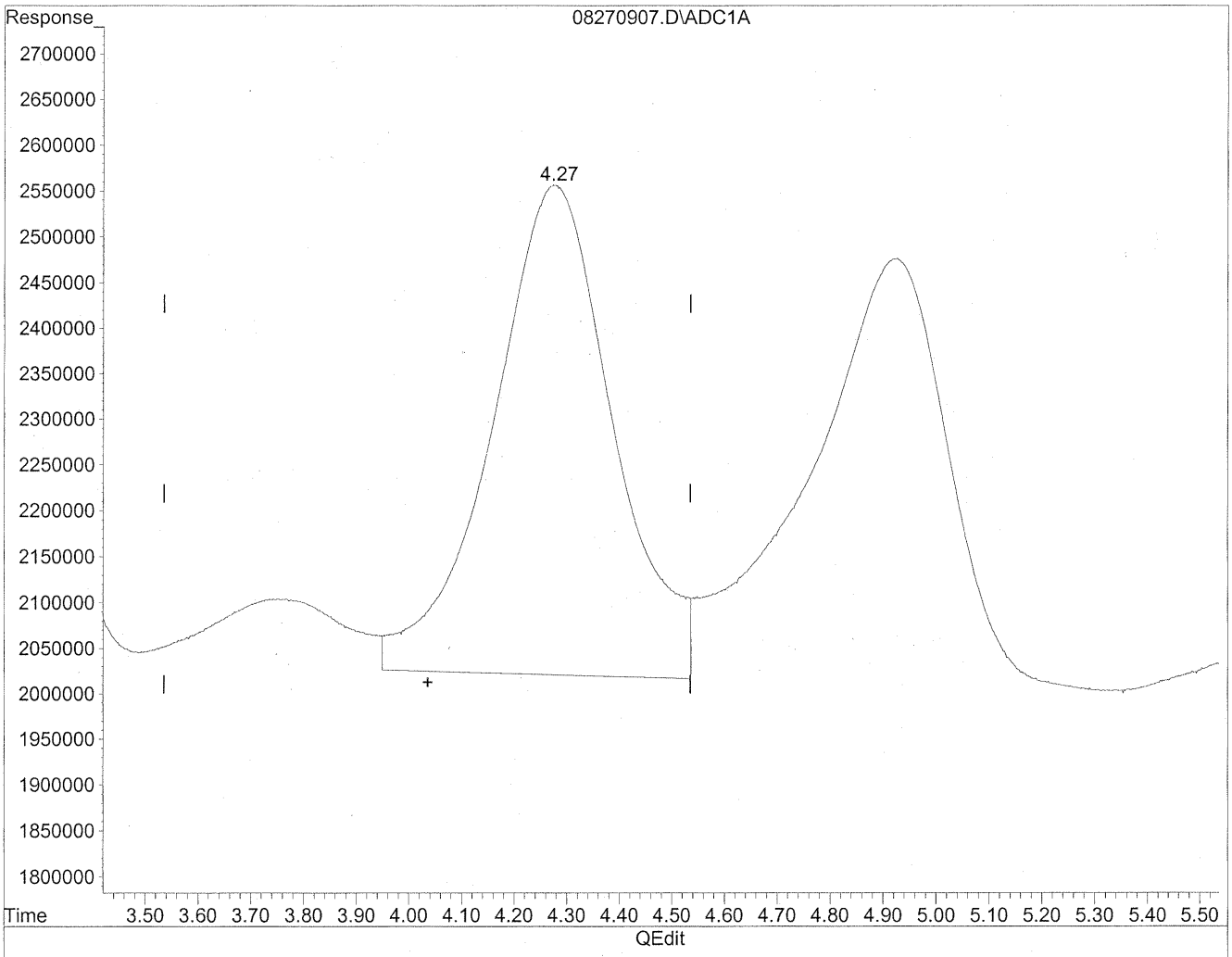
(+) = Expected Retention Time

08270907.D TO110709.M Sun Aug 30 13:33:43 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.28min 875.527ng/ml

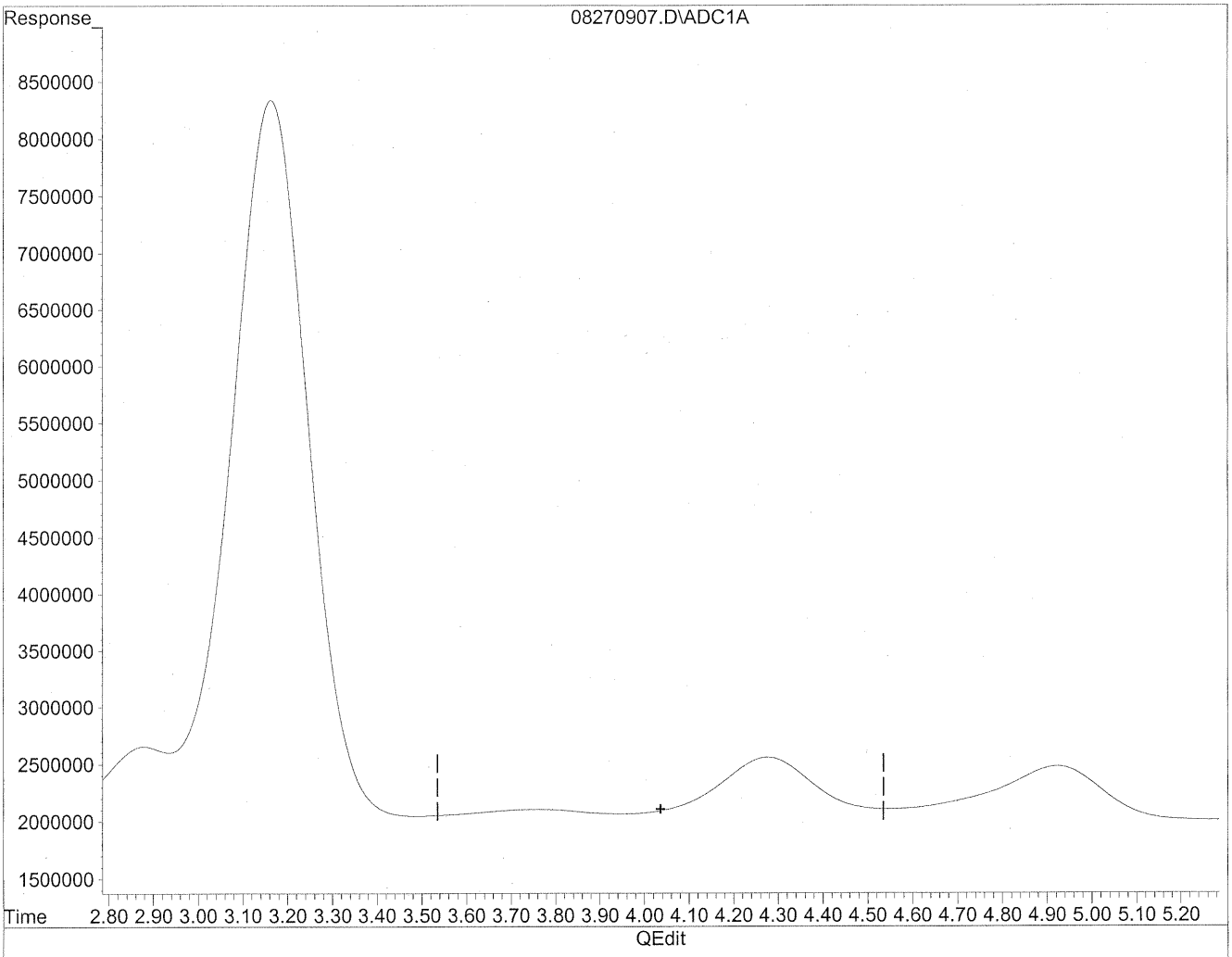
response 85289633

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/31/09 WP

8/31/09

92

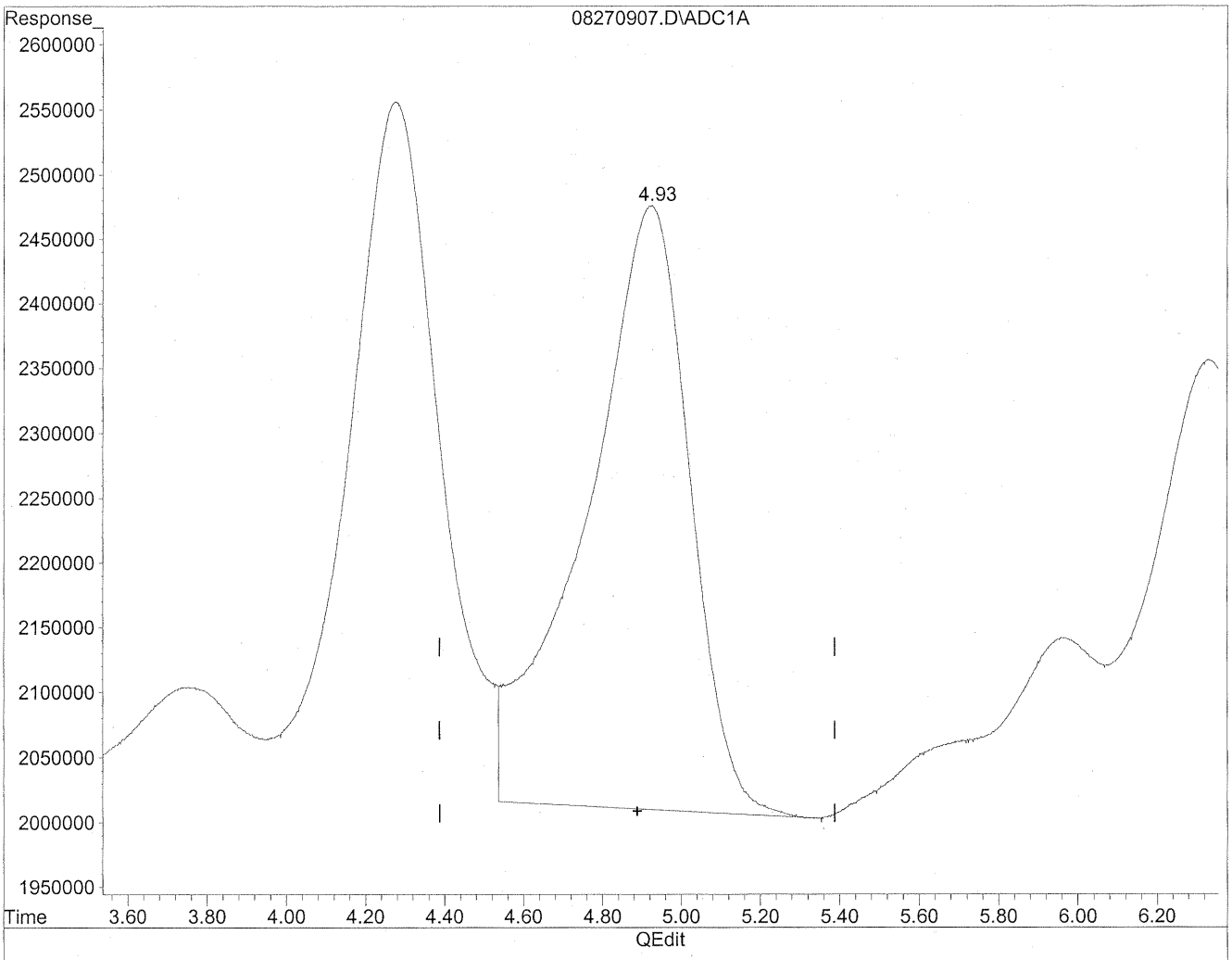
(+) = Expected Retention Time

08270907.D TO110709.M Sun Aug 30 13:33:49 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.92min 962.480ng/ml

response 85021785

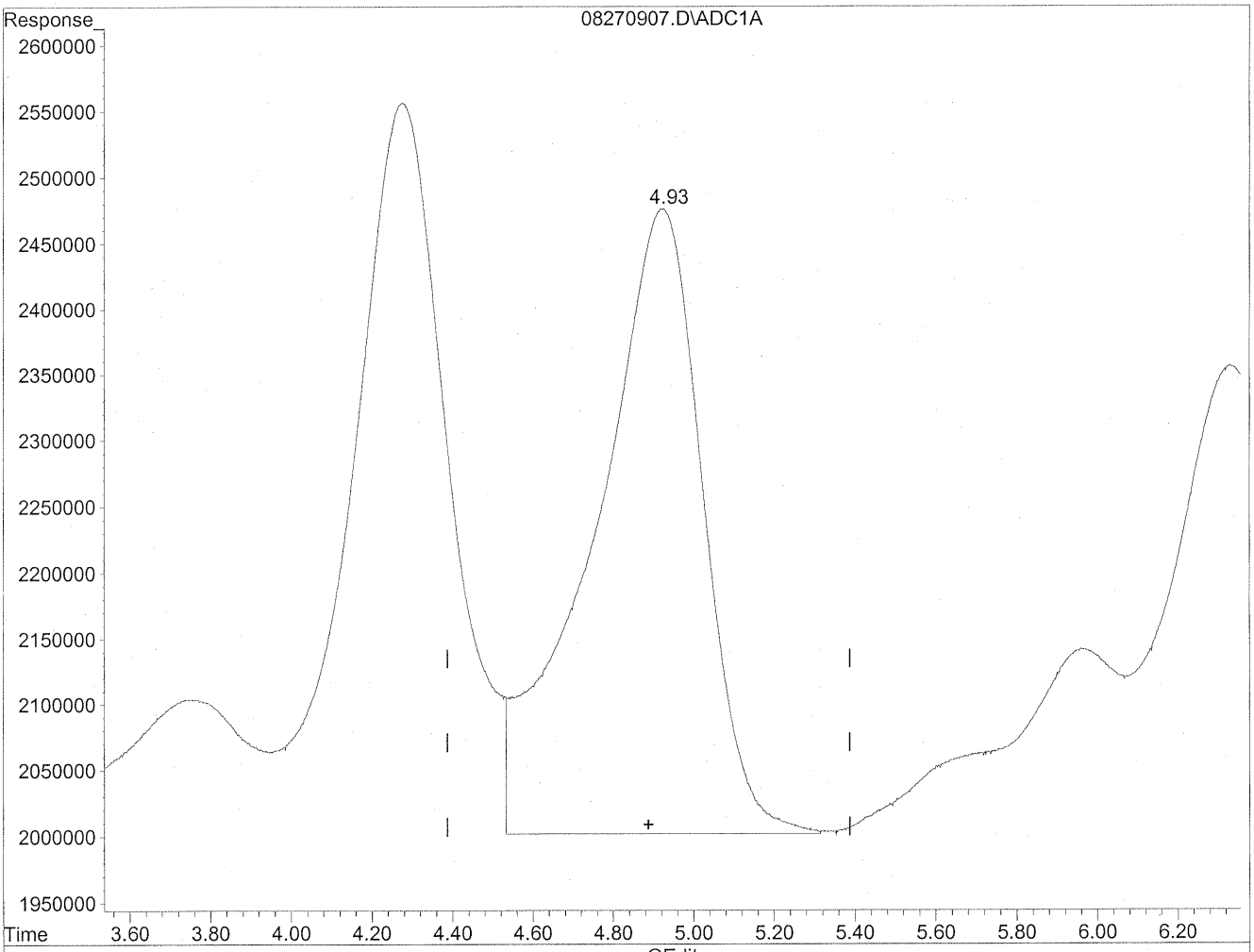
(+) = Expected Retention Time

08270907.D TO110709.M Sun Aug 30 13:33:52 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.93min 1007.640ng/ml m

response 89011049

HC 8/31/09
DC

LW 8/31/09

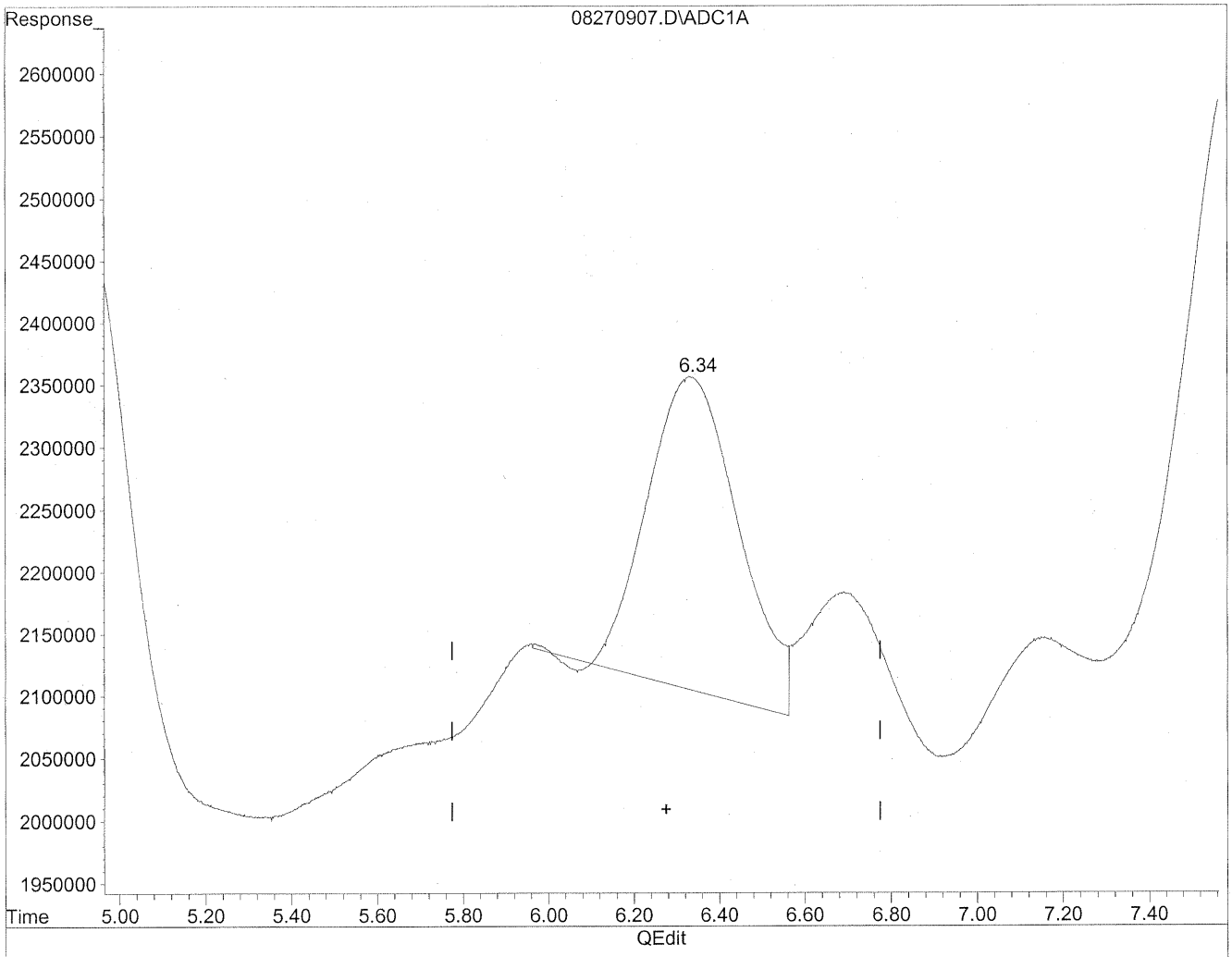
94

(+) = Expected Retention Time
08270907.D TO110709.M Sun Aug 30 13:34:08 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.33min 570.231ng/ml

response 37560690

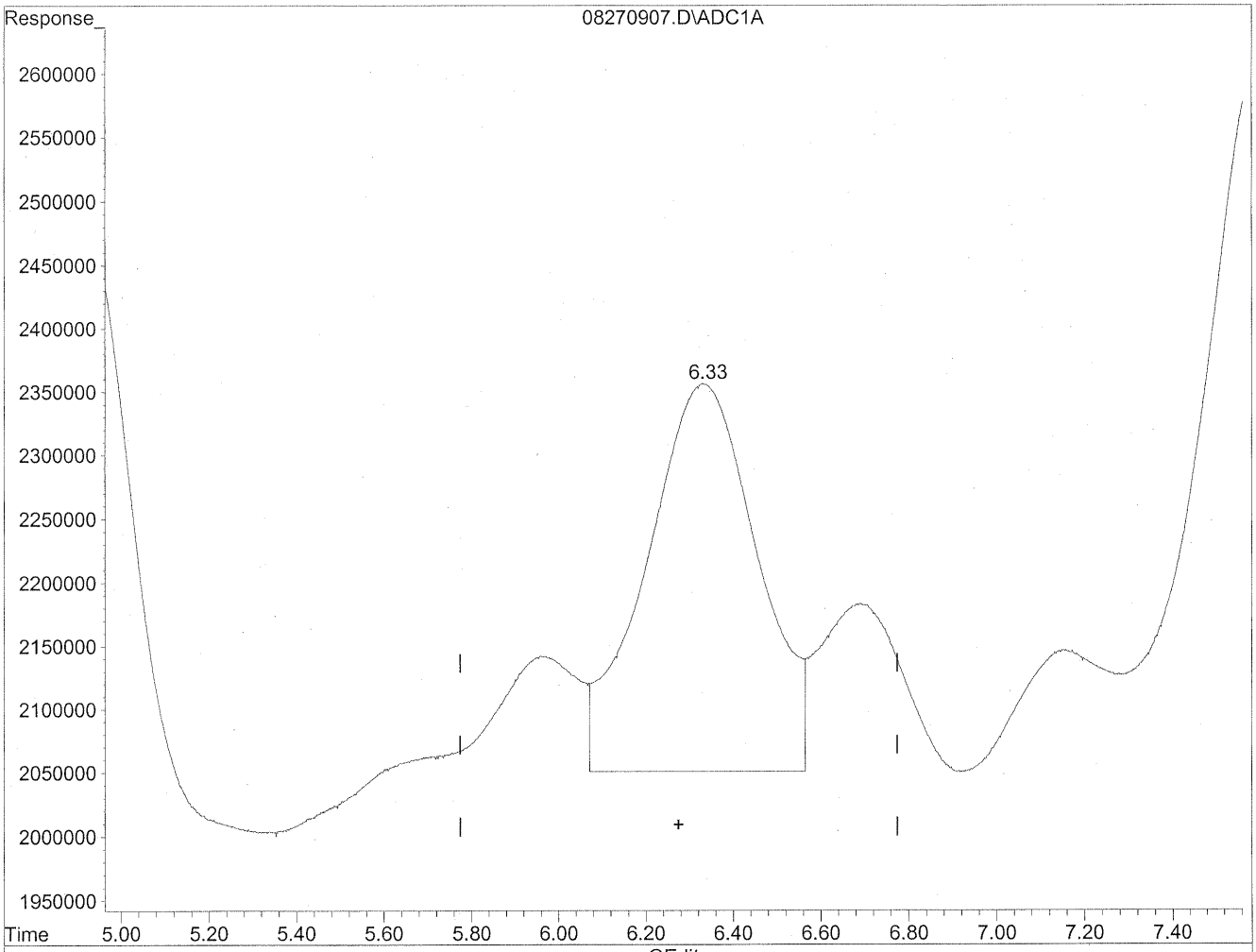
(+) = Expected Retention Time

08270907.D TO110709.M Sun Aug 30 13:34:11 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.33min 820.072ng/ml m

response 54017560

HC 8/31/09 VBL

8/31/09

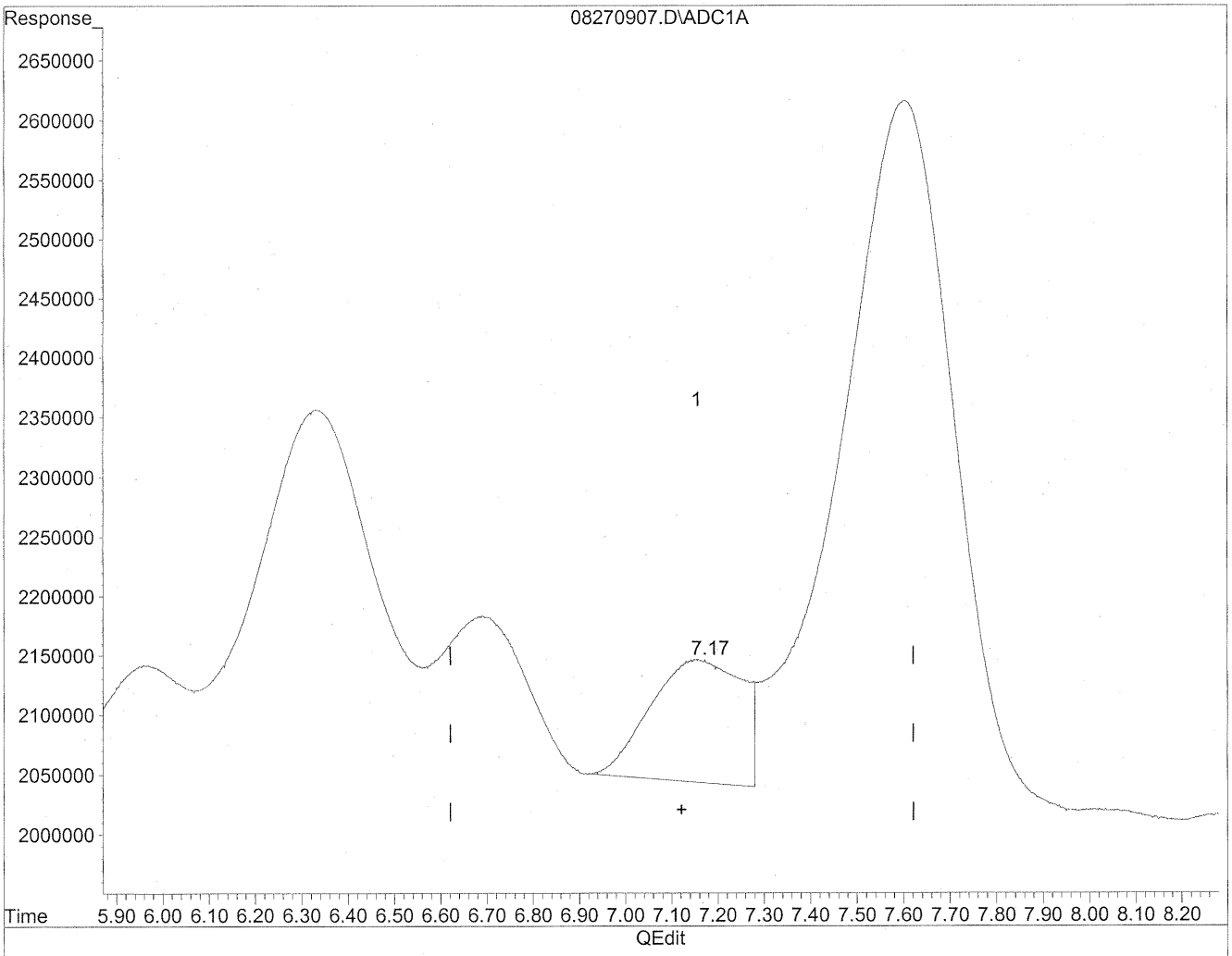
(+) = Expected Retention Time

08270907.D TO110709.M Sun Aug 30 13:34:42 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.15min 180.679ng/ml

response 14138332

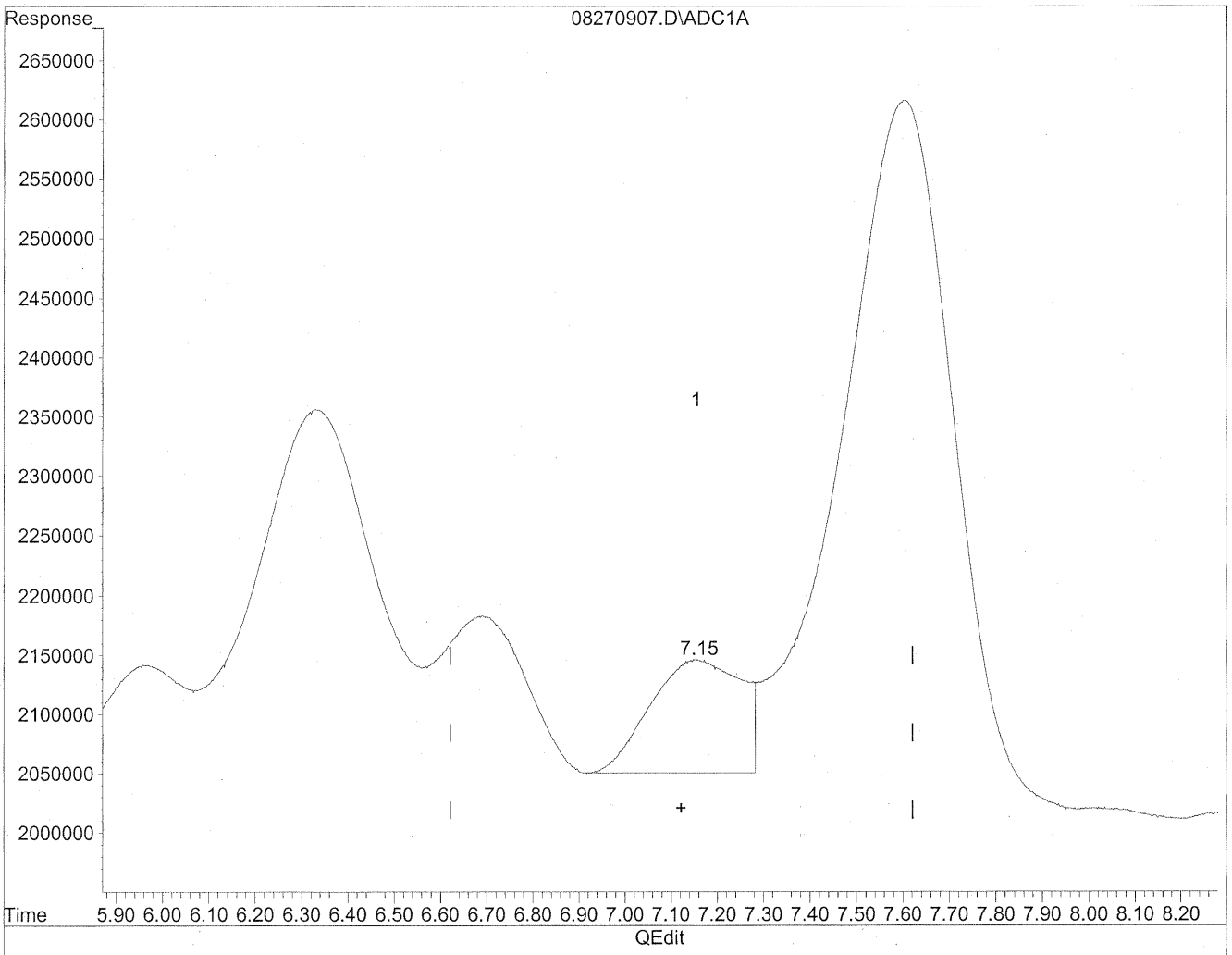
(+) = Expected Retention Time

08270907.D TO110709.M Sun Aug 30 13:34:45 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.15min 167.665ng/ml m

response 13119955

8/31/09 VSC

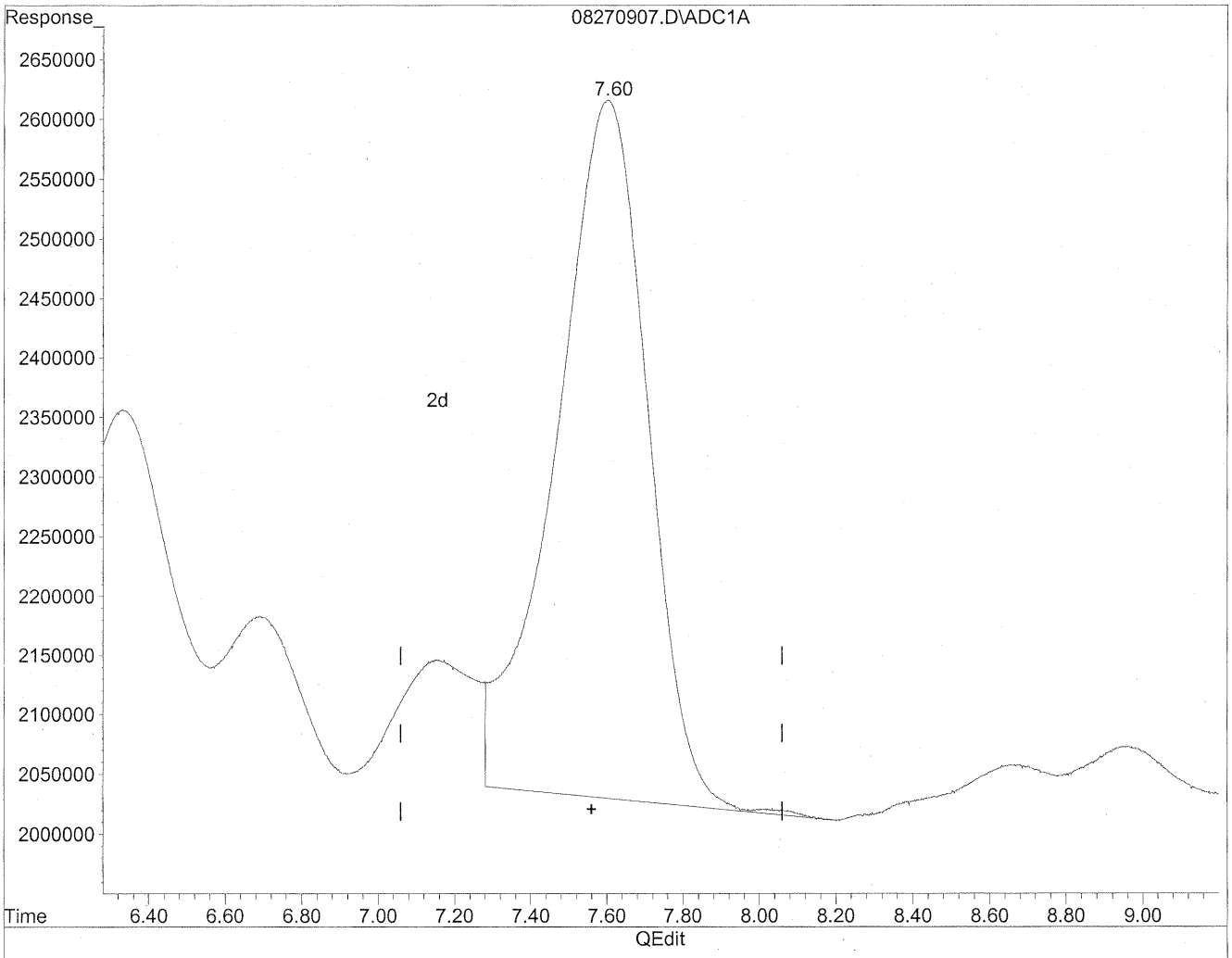
8/31/09

(+) = Expected Retention Time
08270907.D TO110709.M Sun Aug 30 13:34:56 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.60min 1338.505ng/ml

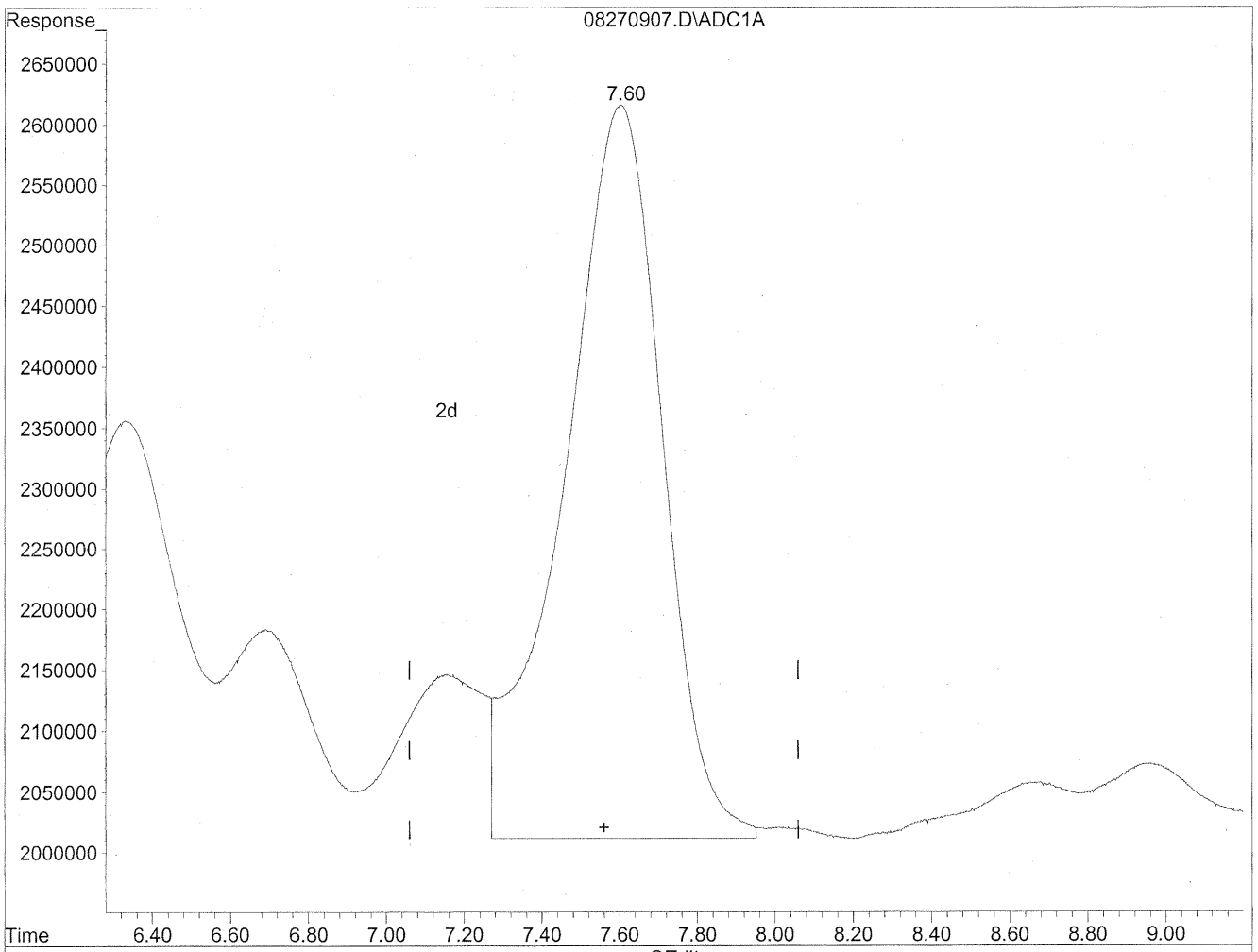
response 98386817

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.60min 1440.994ng/ml m

response 105920278

8/31/09

8/31/09

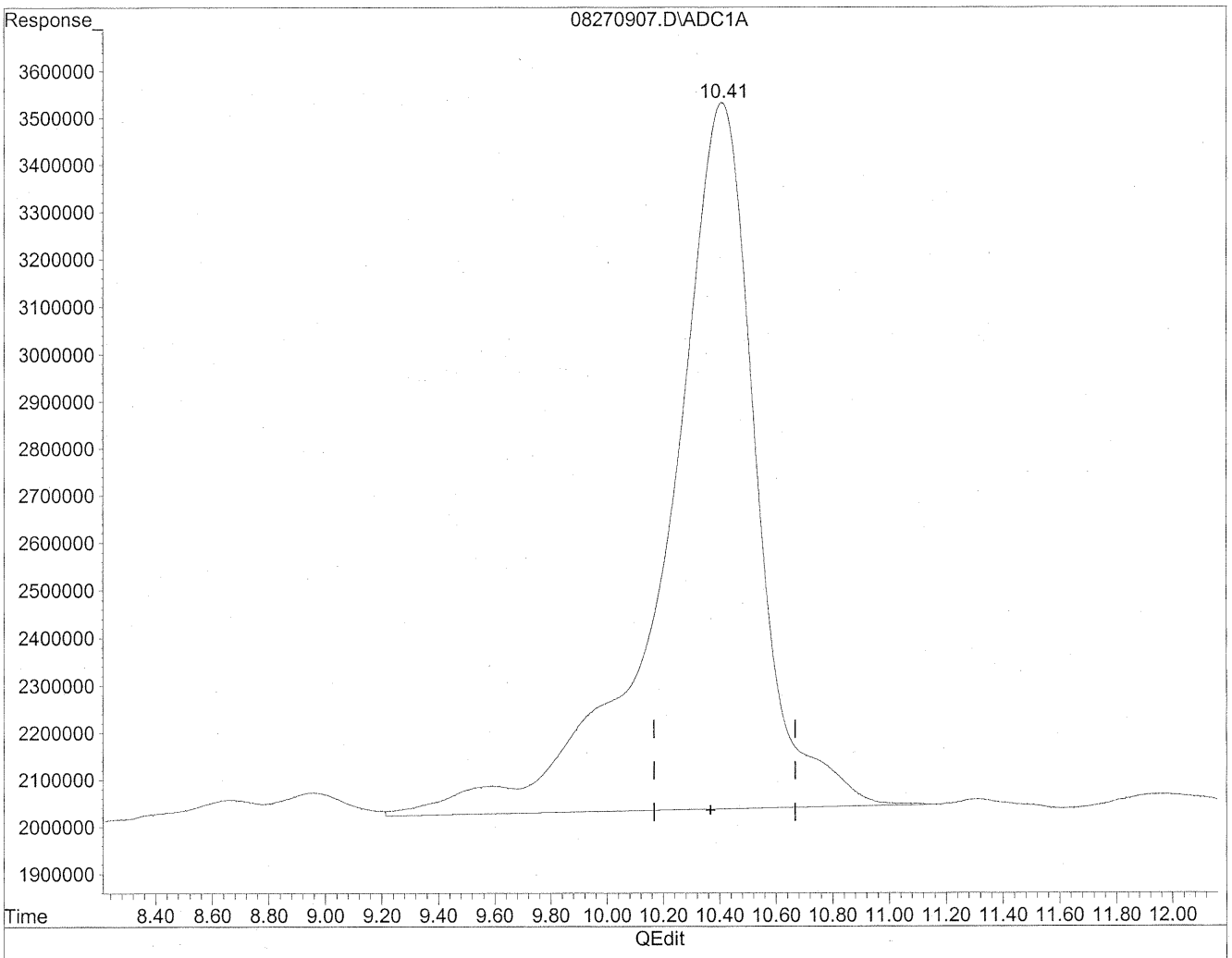
100

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.41min 4937.092ng/ml

response 332482325

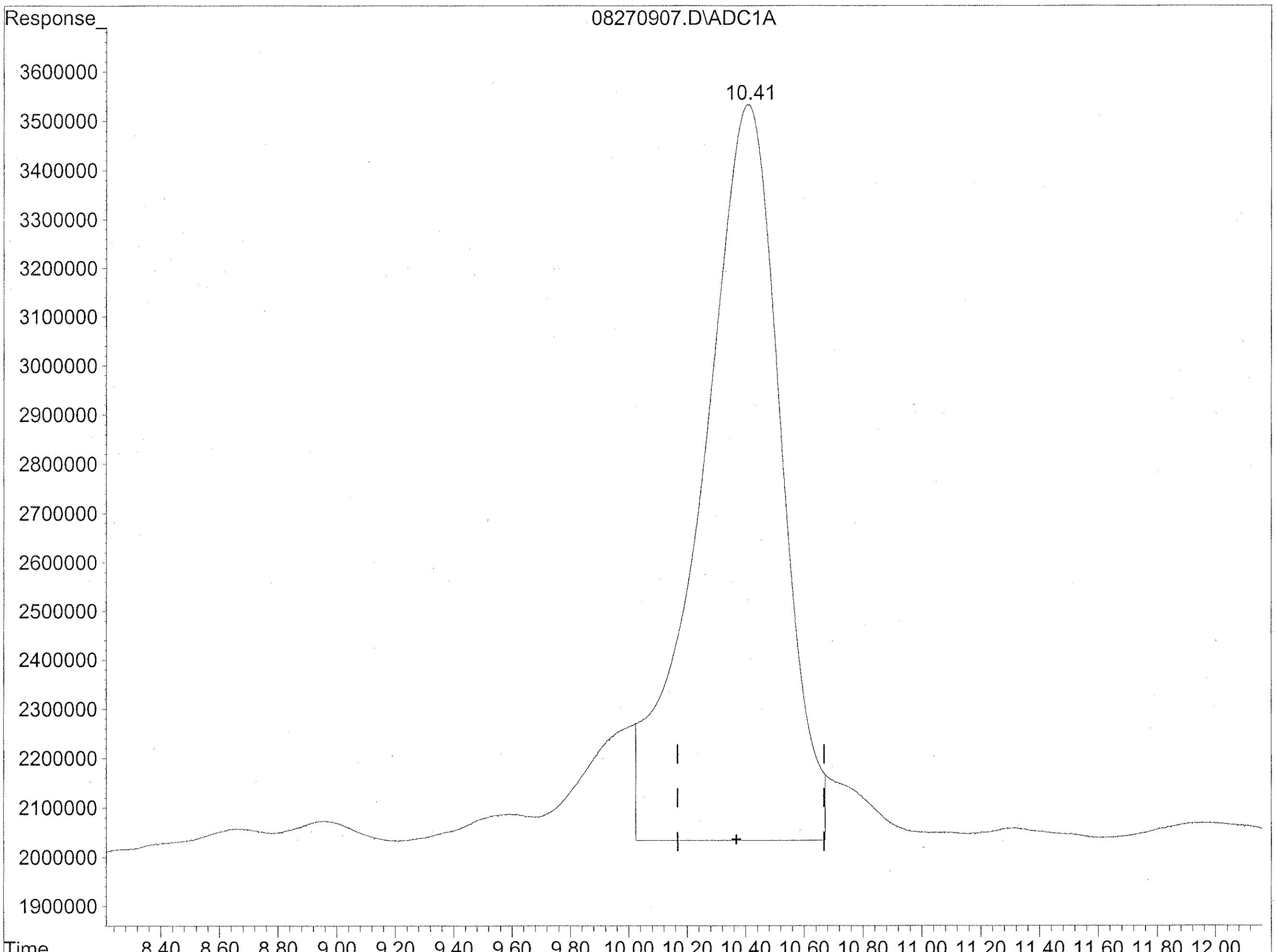
101

(+) = Expected Retention Time

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D      Vial: 7
Acq On    : 27 Aug 2009  10:35 am                      Operator: HC
Sample    : P0902942-004 front 1.0ml                   Inst    : LC 01
Misc      :                                            Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Aug 30 13:33 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   : Sat Aug 29 17:49:00 2009
Response via  : Multiple Level Calibration
```



QEdit

(11) Hexaldehyde

10.41min 4193.832ng/ml m

response 282428395

HC 8/31/09 SH IBC

WP 8/31/09

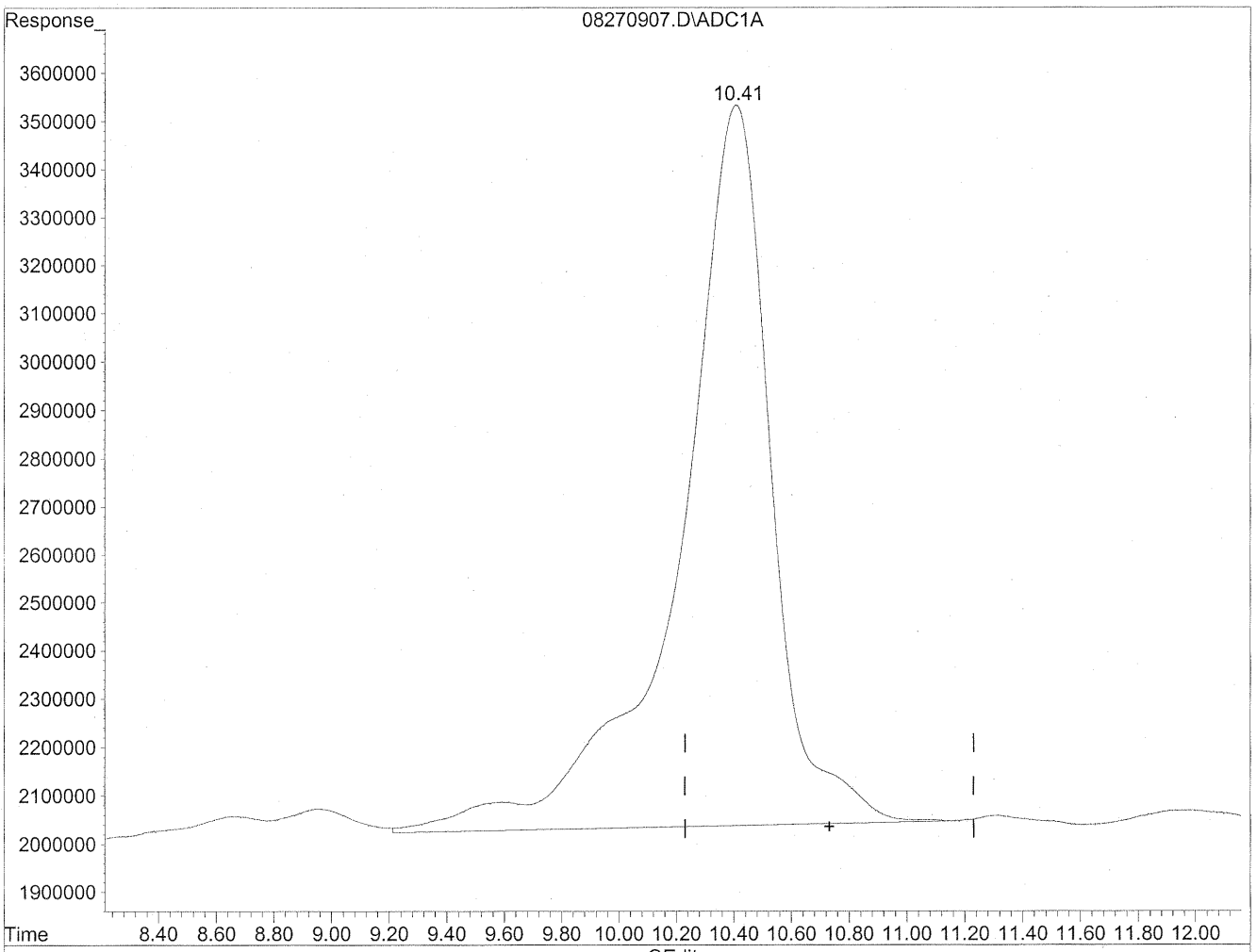
(+) = Expected Retention Time

08270907.D TO110709.M Sun Aug 30 13:35:59 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.41min 6783.500ng/ml

response 332482325

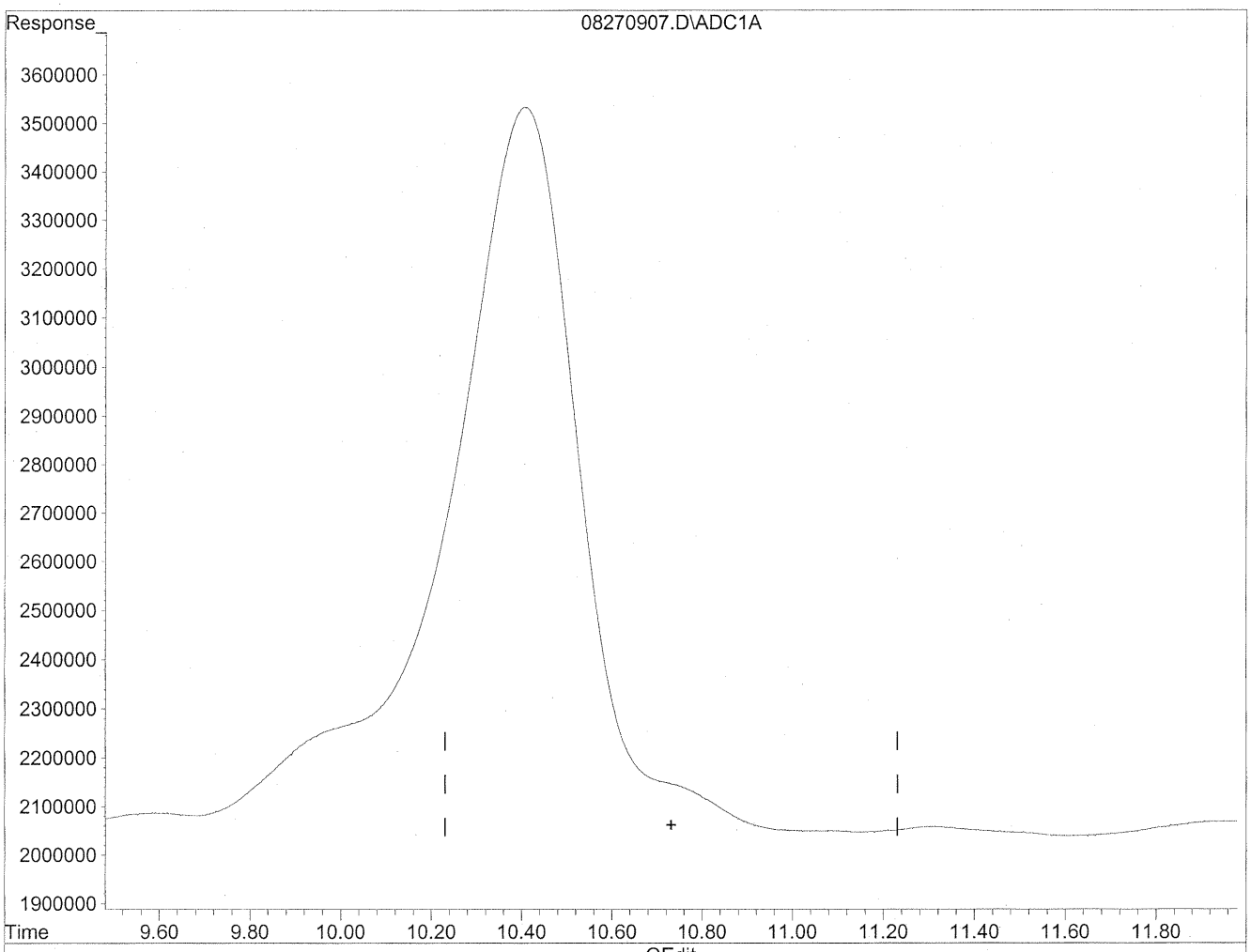
(+) = Expected Retention Time

08270907.D TO110709.M Sun Aug 30 13:36:02 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270907.D Vial: 7
Acq On : 27 Aug 2009 10:35 am Operator: HC
Sample : P0902942-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:33 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 : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/31/09 MP

8/31/09

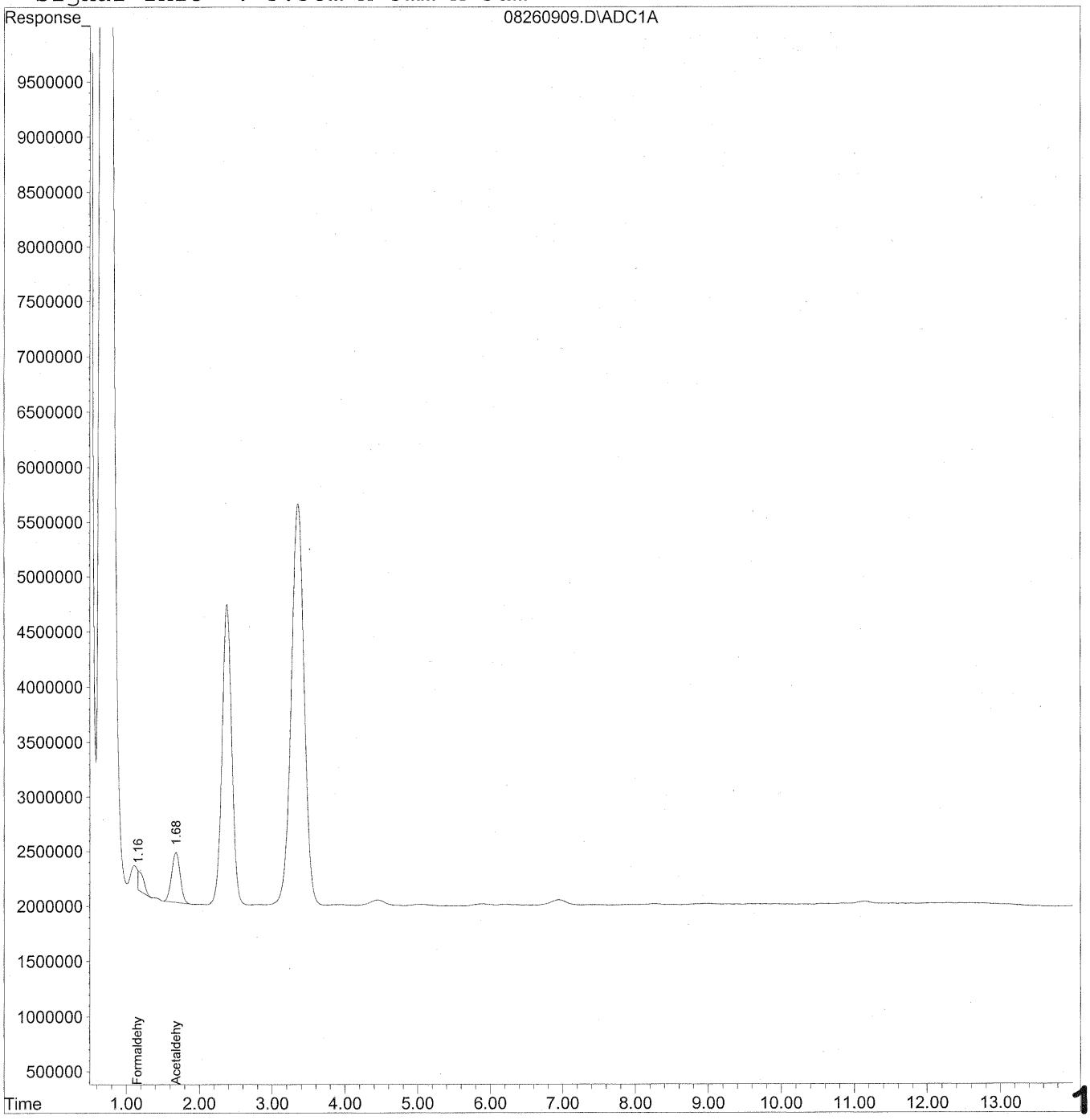
104

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_08\26\08260909.D Vial: 9
Acq On : 26 Aug 2009 7:05 pm Operator: HC
Sample : P0902942-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:49 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 : Thu Aug 27 07:35:56 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



105

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\26\08260909.D Vial: 9
Acq On : 26 Aug 2009 7:05 pm Operator: HC
Sample : P0902942-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:49 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 : Thu Aug 27 07:35:56 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

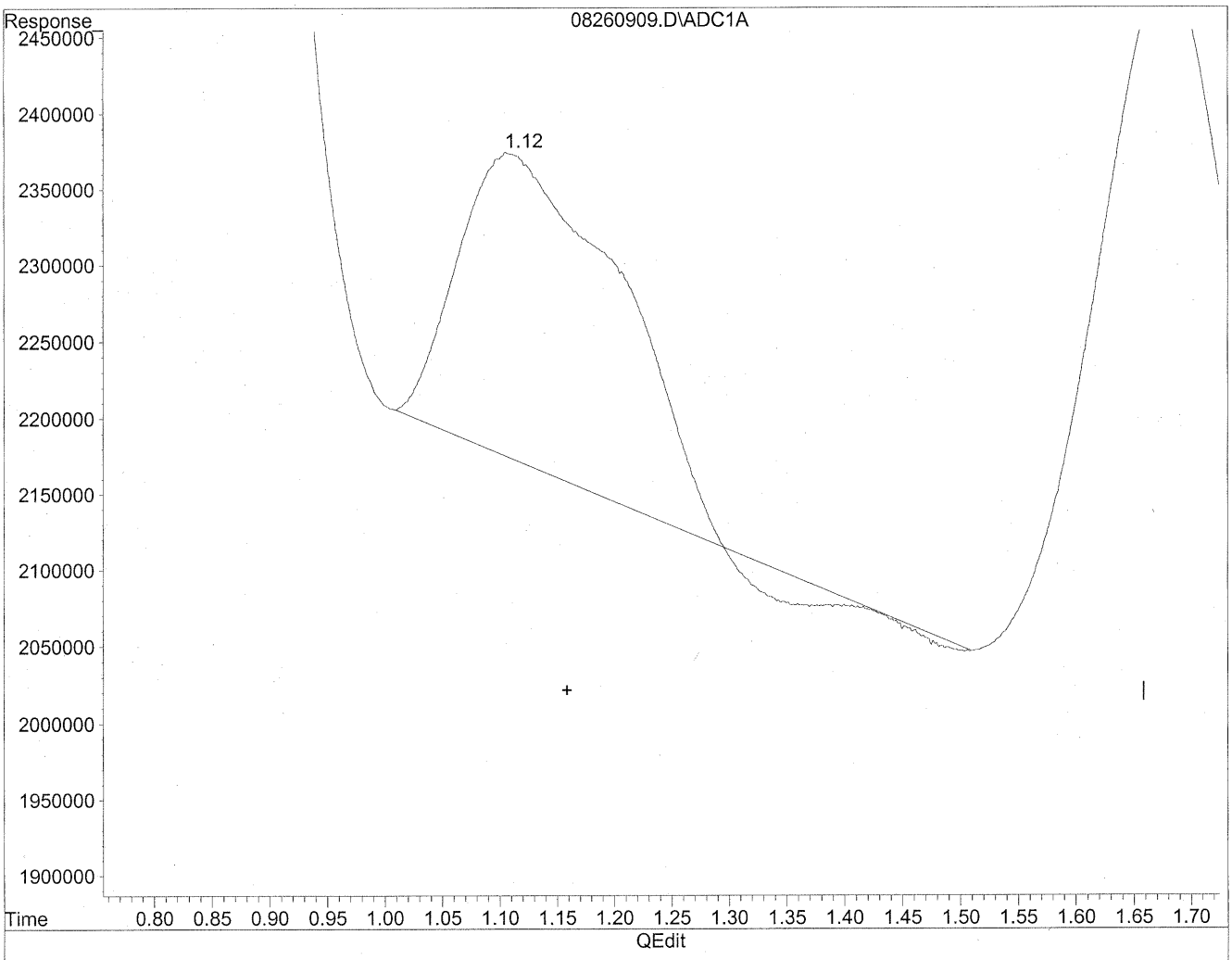
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	9763540	53.184	ng/mlm
2)	Acetaldehyde	1.68	37908109	270.341	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	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\26\08260909.D Vial: 9
Acq On : 26 Aug 2009 7:05 pm Operator: HC
Sample : P0902942-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.11min 105.450ng/ml

response 19358588

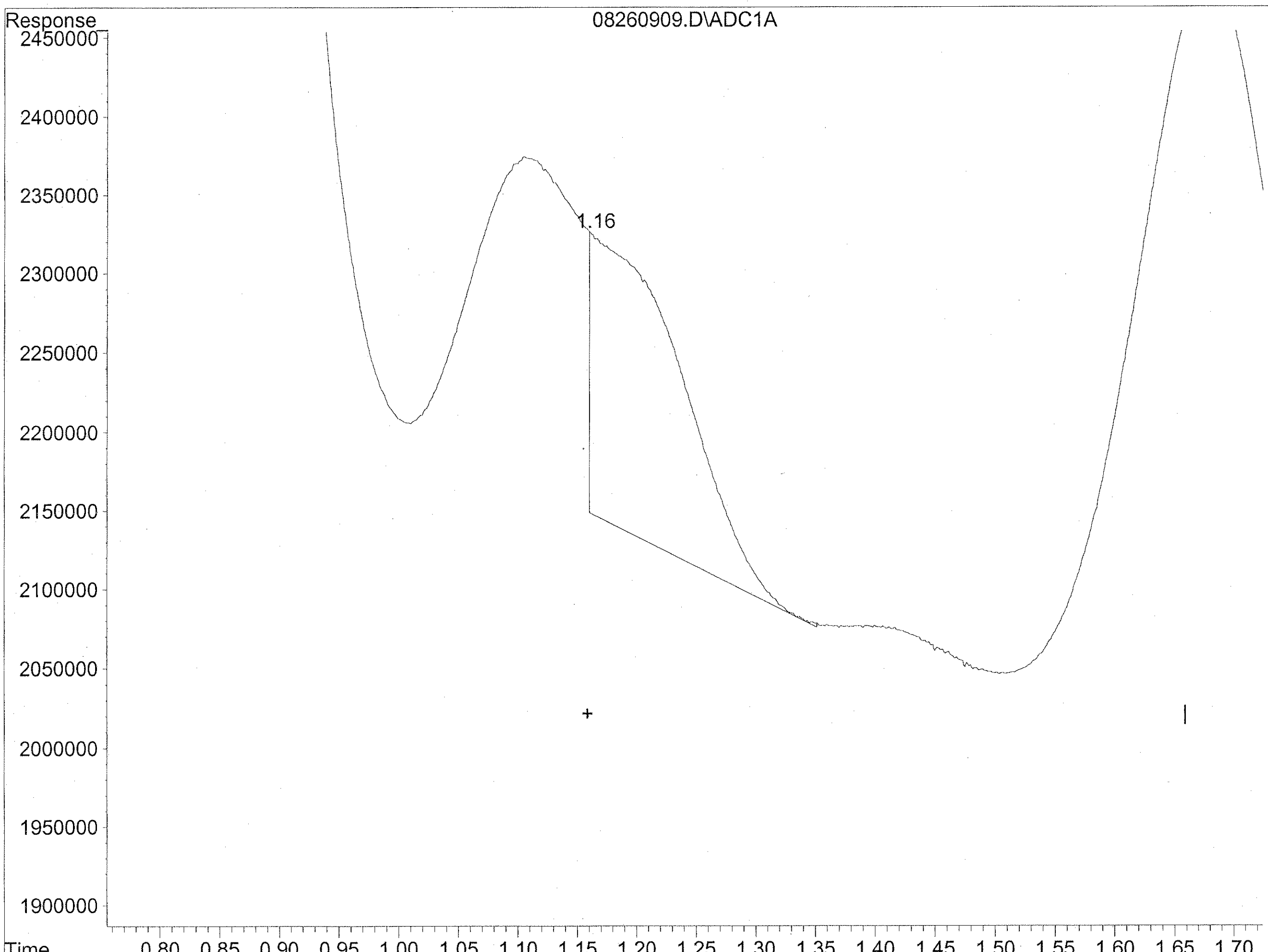
(+) = Expected Retention Time

08260909.D TO110709.M Sat Aug 29 16:49:07 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260909.D   Vial: 9
Acq On   : 26 Aug 2009   7:05 pm   Operator: HC
Sample   : P0902942-004 back 1.0ml   Inst   : LC 01
Misc   :   Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method   : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title   : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.16min 53.184ng/ml m

response 9763540

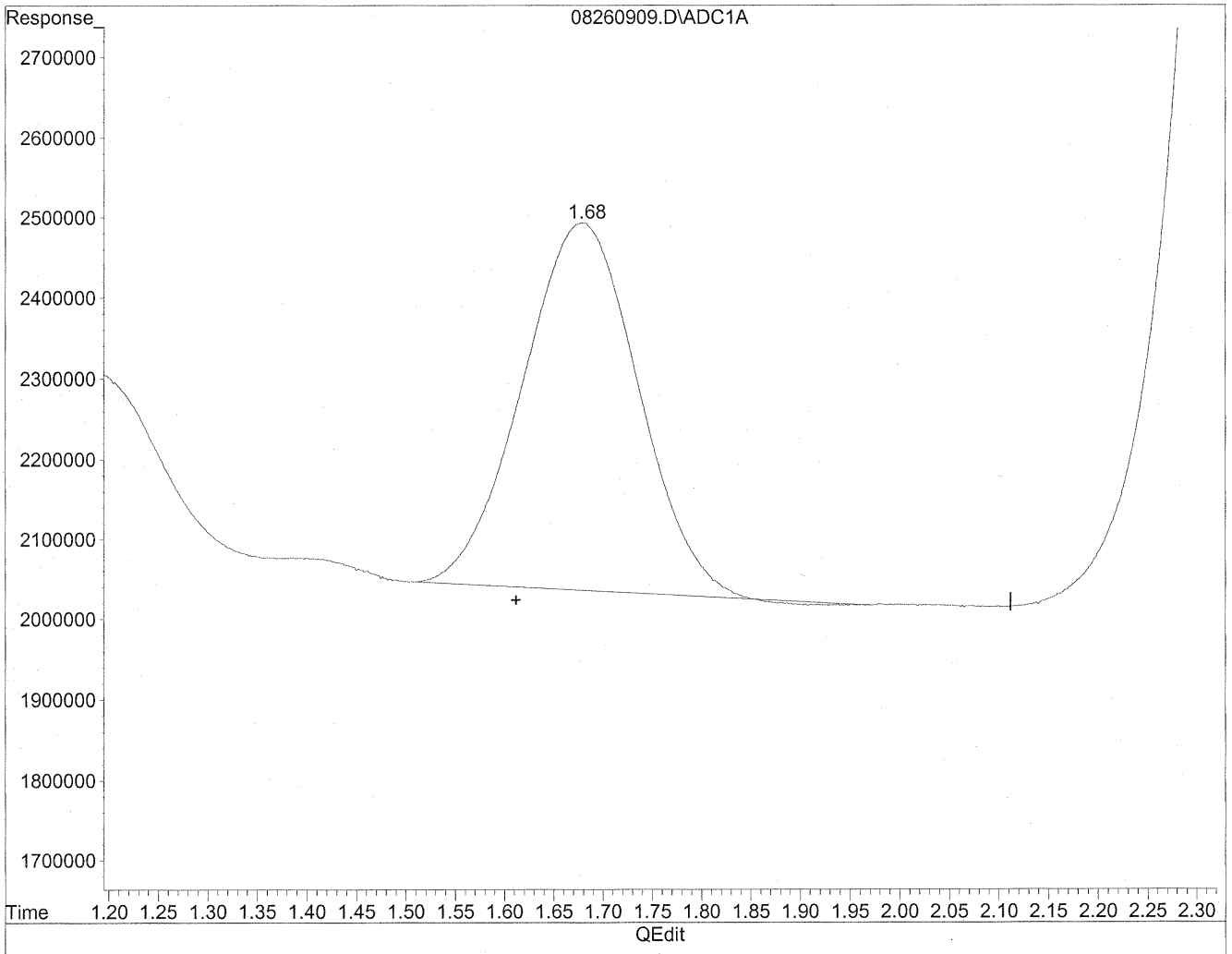
HC 8/30/09
SP
[illegible] 8/31/09

(+) = Expected Retention Time
08260909.D TO110709.M   Sat Aug 29 16:49:24 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260909.D Vial: 9
Acq On : 26 Aug 2009 7:05 pm Operator: HC
Sample : P0902942-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.68min 266.426ng/ml

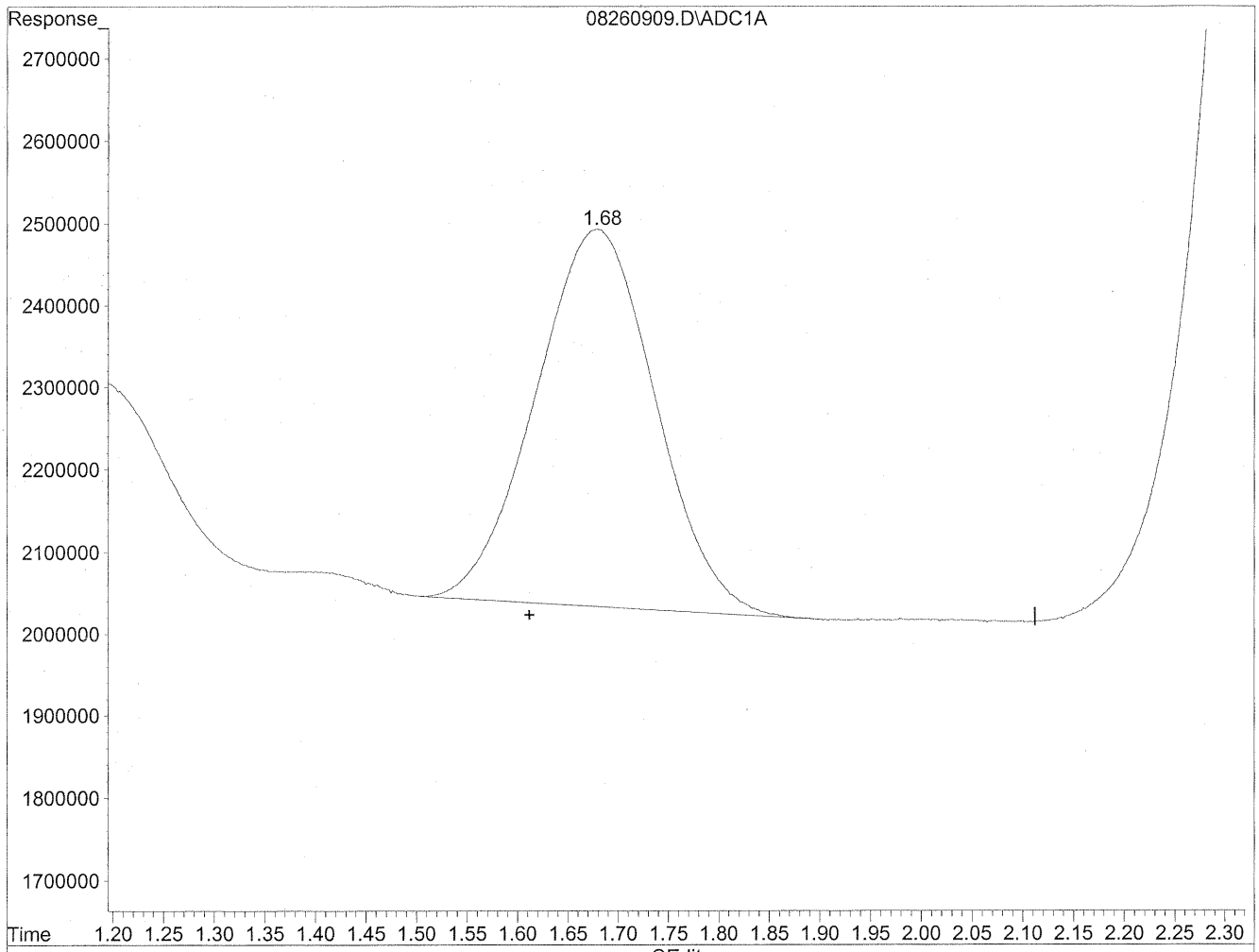
response 37359204

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260909.D Vial: 9
Acq On : 26 Aug 2009 7:05 pm Operator: HC
Sample : P0902942-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.68min 270.341ng/ml m

response 37908109

8/31/09

8/31/09

110

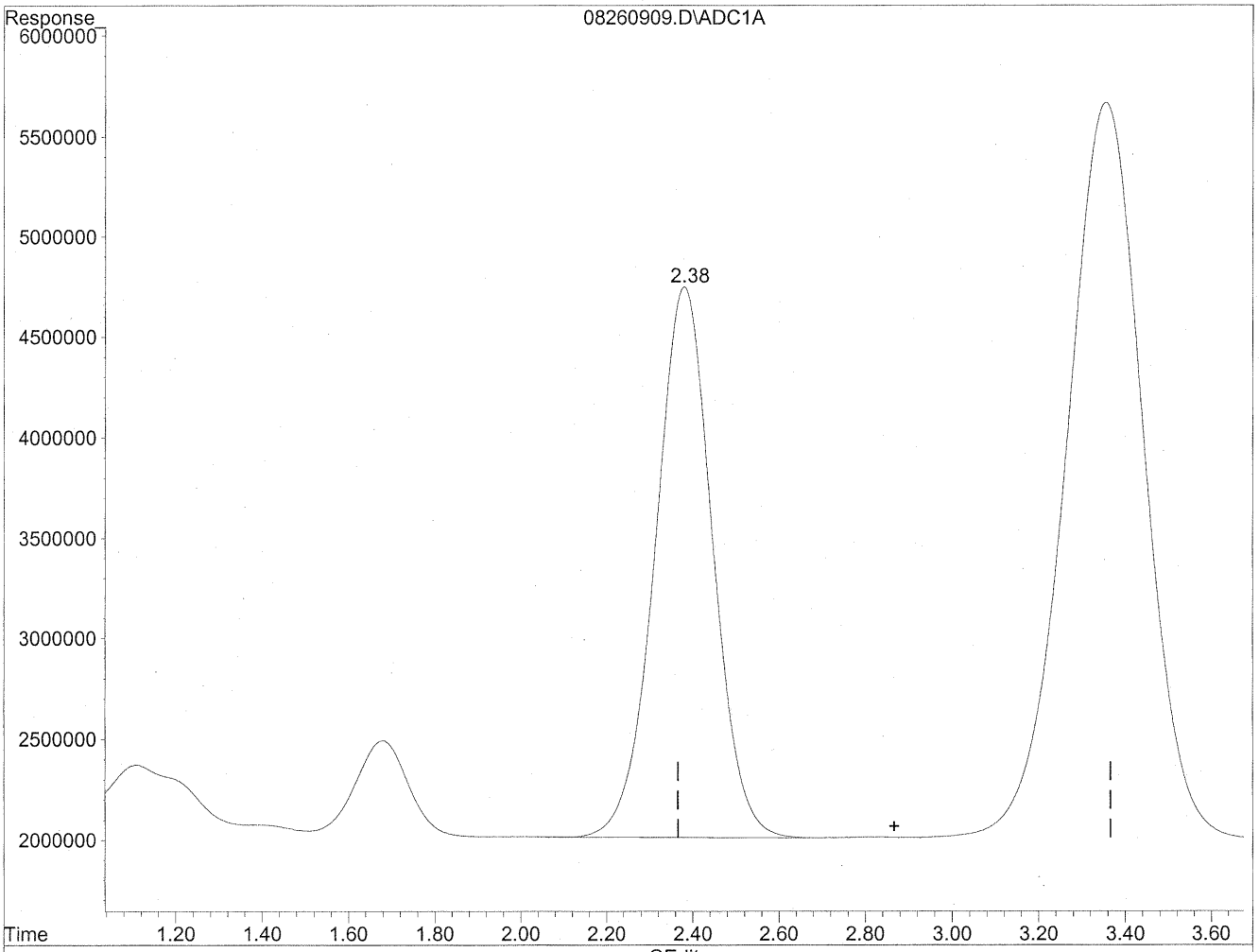
(+) = Expected Retention Time

08260909.D TO110709.M Sat Aug 29 16:49:40 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260909.D Vial: 9
Acq On : 26 Aug 2009 7:05 pm Operator: HC
Sample : P0902942-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.38min 2360.815ng/ml

response 251887700

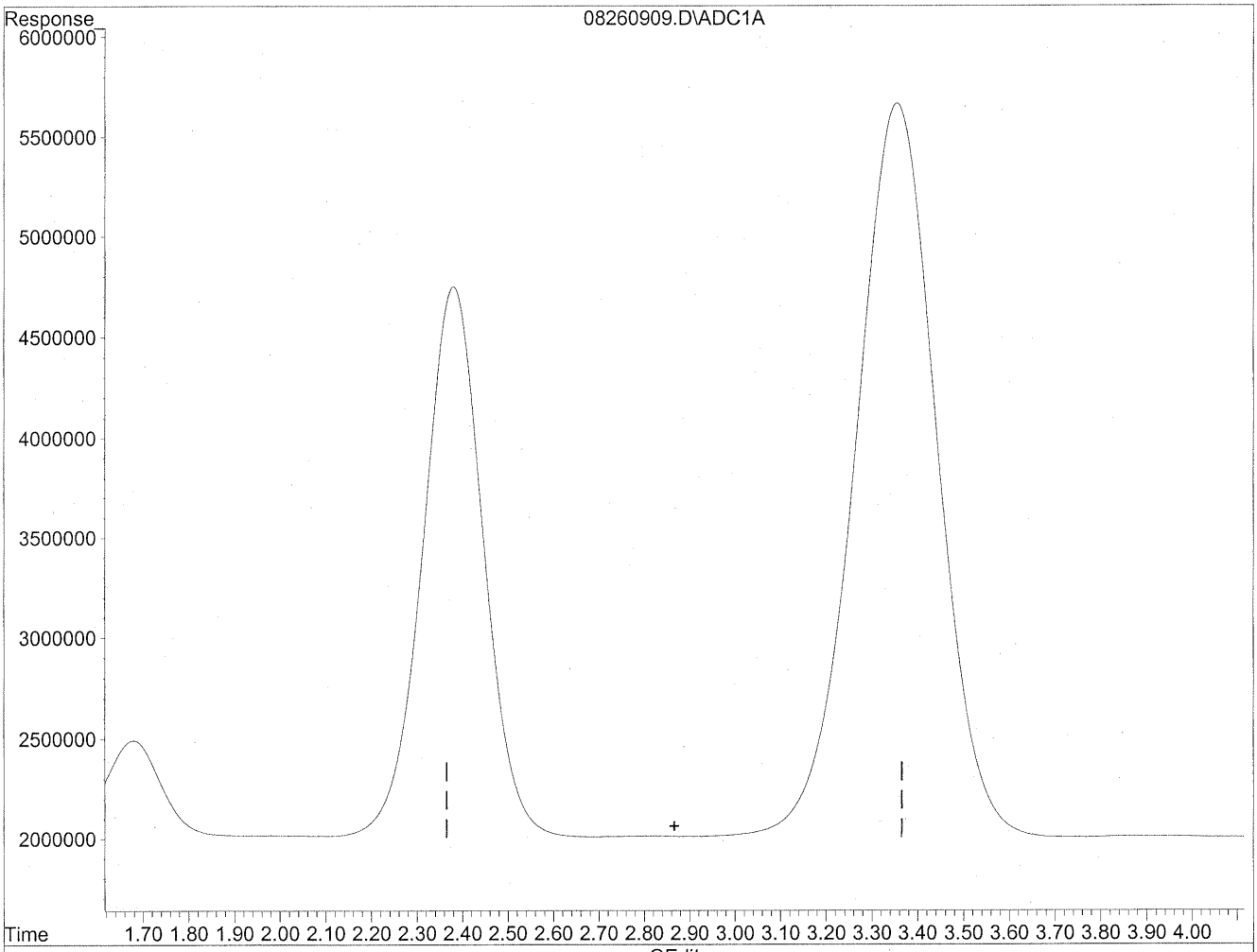
111

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260909.D Vial: 9
Acq On : 26 Aug 2009 7:05 pm Operator: HC
Sample : P0902942-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/30/09 MP

8/31/09

112

(+) = Expected Retention Time

08260909.D TO110709.M Sat Aug 29 16:49:47 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **101473**
Client Project ID: **16512**

CAS Project ID: P0902942
CAS Sample ID: P0902942-005

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: 8/24/09
Date Received: 8/25/09
Date Analyzed: 8/26-27/09
Desorption Volume: 1.0 ml
Volume Sampled: 104.54 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**2,300**	**22**	0.96	**18**	0.78	
75-07-0	**Acetaldehyde**	**4,800**	**46**	0.96	**25**	0.53	
123-38-6	**Propionaldehyde**	**640**	**6.1**	0.96	**2.6**	0.40	
4170-30-3	Crotonaldehyde, Total	< 200	ND	1.9	ND	0.67	**V**
123-72-8	**Butyraldehyde**	**1,100**	**11**	0.96	**3.6**	0.32	**M**
100-52-7	**Benzaldehyde**	**800**	**7.6**	0.96	**1.8**	0.22	
590-86-3	**Isovaleraldehyde**	**140**	**1.3**	0.96	**0.37**	0.27	
110-62-3	**Valeraldehyde**	**1,400**	**14**	0.96	**3.9**	0.27	
529-20-4	o-Tolualdehyde	< 100	ND	0.96	ND	0.19	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	1.9	ND	0.39	
66-25-1	**n-Hexaldehyde**	**5,000**	**48**	0.96	**12**	0.23	
5779-94-2	2,5-Dimethylbenzaldehyde	< 200	ND	1.9	ND	0.35	**V**

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.
M = Matrix interference; results may be biased high.
V = The continuing calibration verification standard was outside (biased low) the specified limits for this compound.

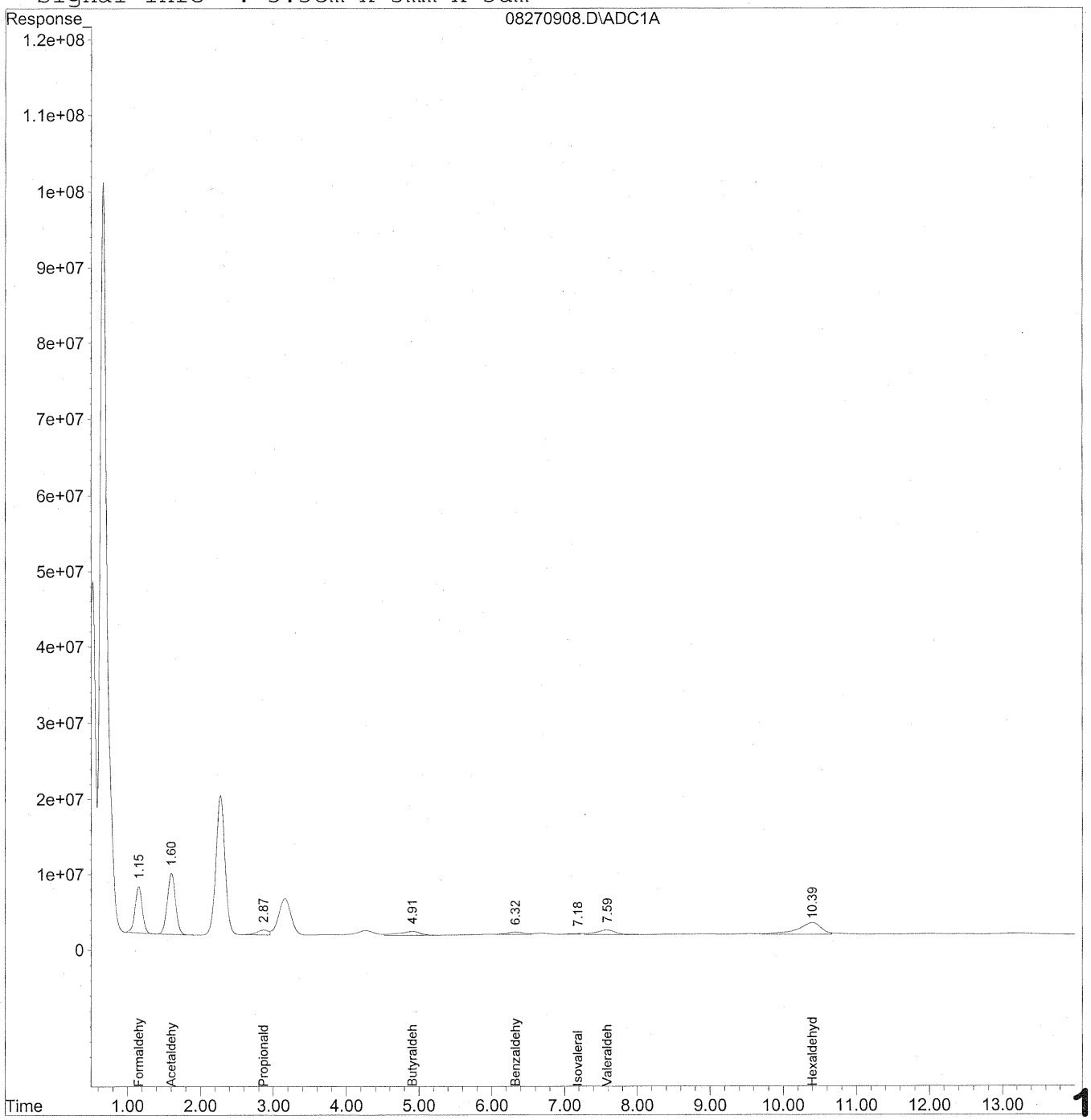
Verified By: Re Date: 9/11/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:39 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 : Thu Aug 27 08:33:51 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\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:39 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 : Thu Aug 27 08:33:51 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

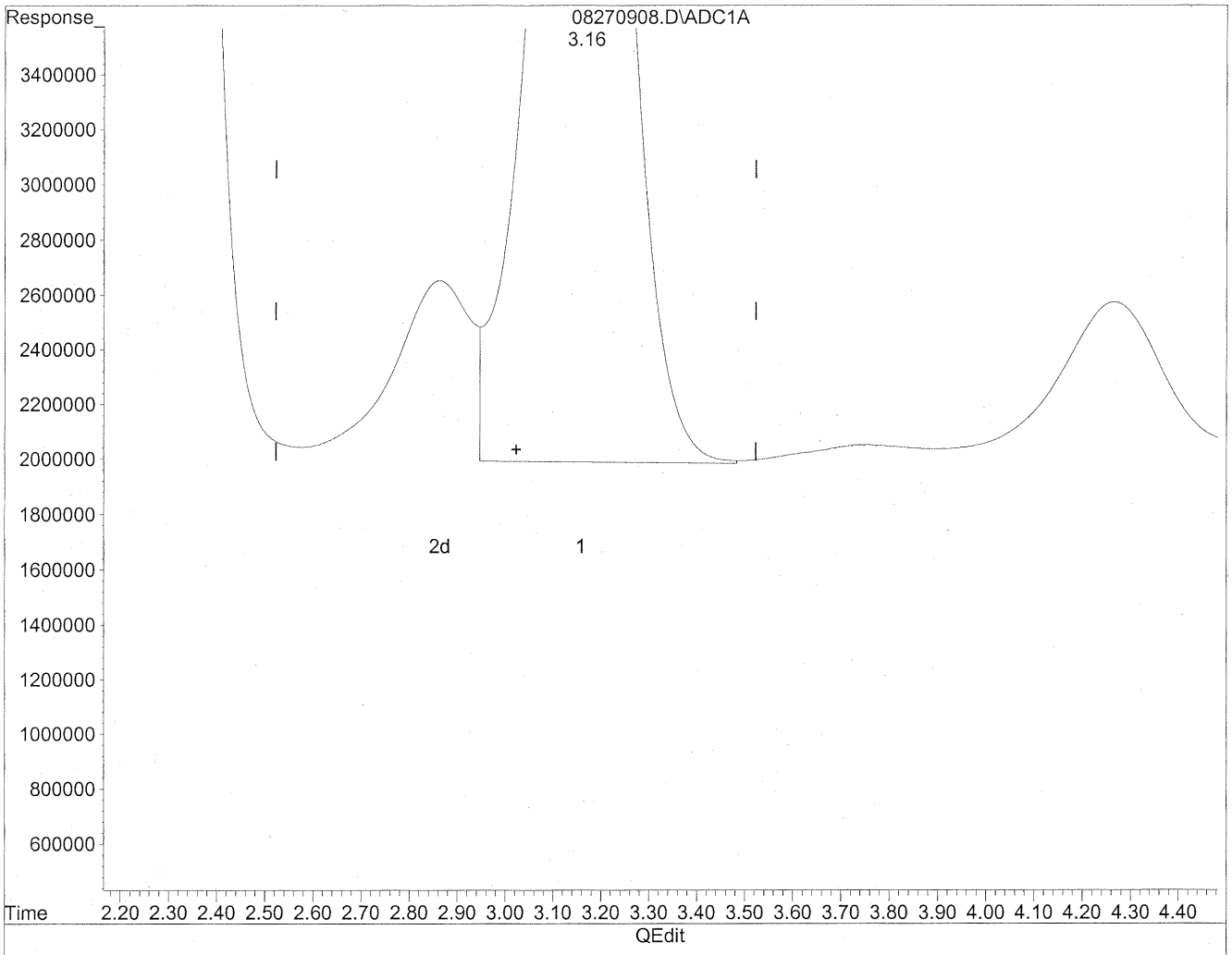
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.15	416529039	2268.907	ng/ml
2)	Acetaldehyde	1.60	635629932	4532.977	ng/ml
3)	Propionaldehyde	2.87	67772106	635.193	ng/mlm
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	4.91	97226838	1100.646	ng/mlm
6)	Benzaldehyde	6.32	52437346	796.082	ng/mlm
7)	Isovaleraldehyde	7.18	10723919	137.045	ng/mlm
8)	Valeraldehyde	7.59	105159475	1430.644	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	10.39f	338148817	5021.234	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\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



QEdit

(3) Propionaldehyde

3.16min 5339.393ng/ml

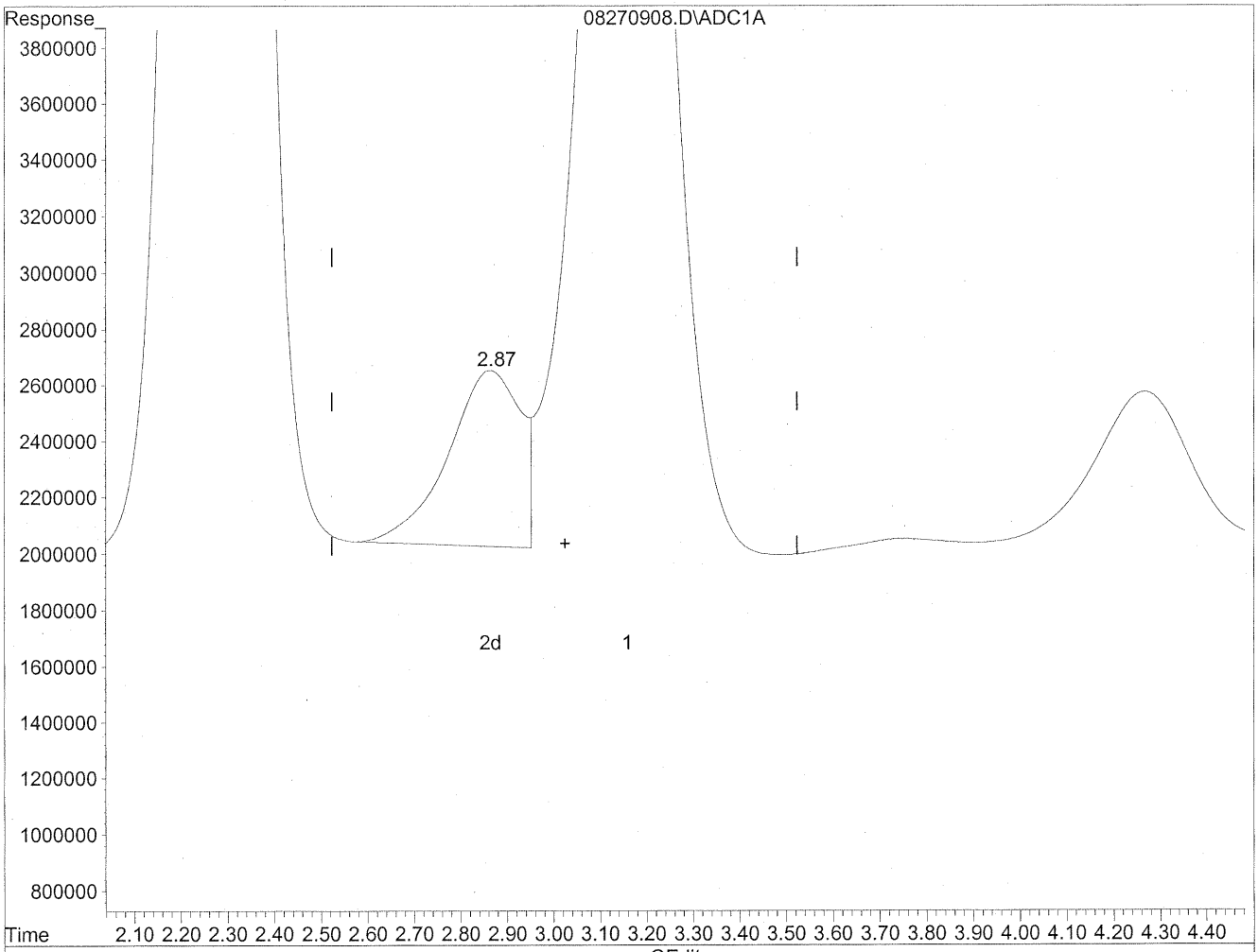
response 569687758

(+) = Expected Retention Time
08270908.D TO110709.M Sun Aug 30 13:37:08 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.87min 635.193ng/ml m

response 67772106

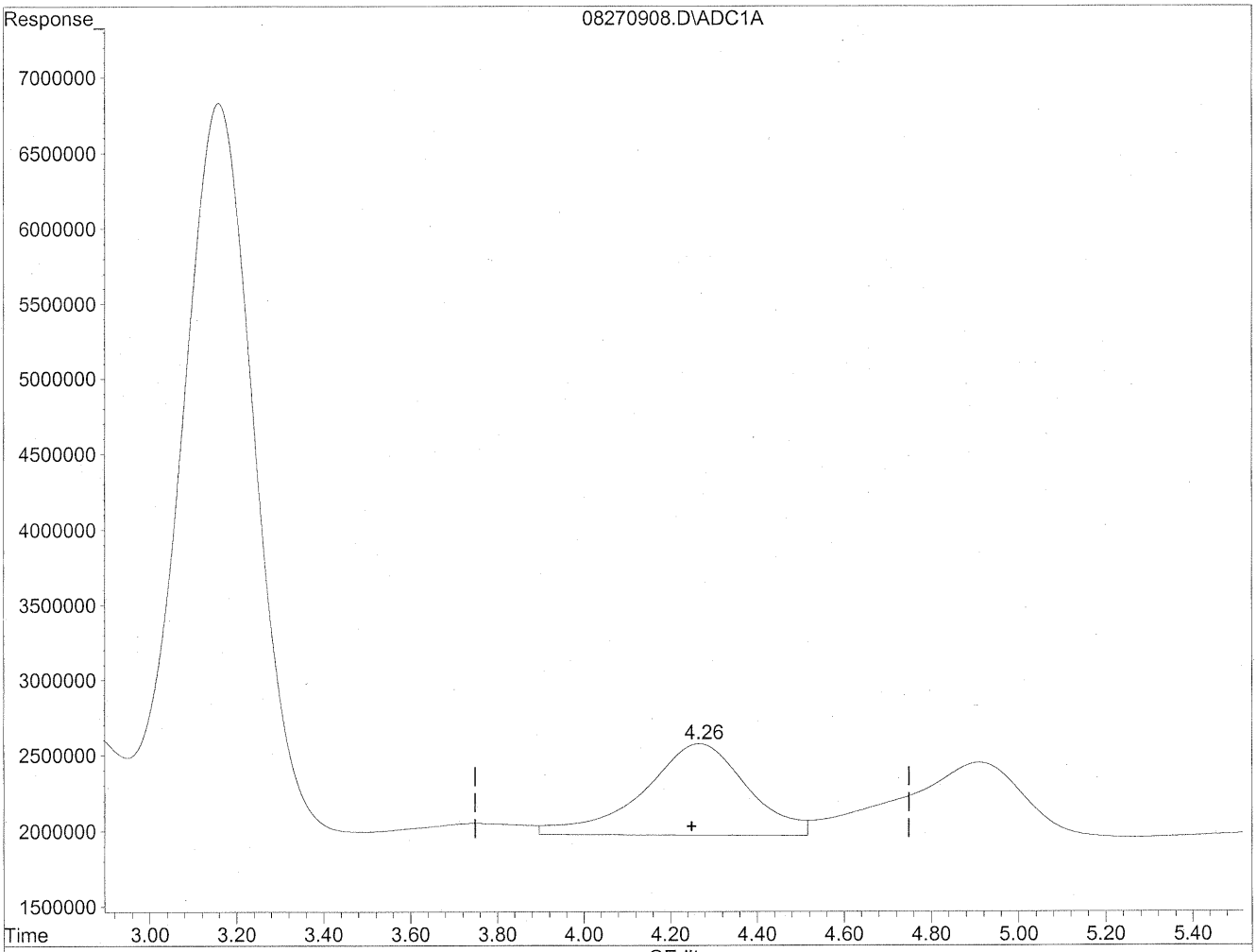
(+) = Expected Retention Time

08270908.D TO110709.M Sun Aug 30 13:37:25 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.27min 1034.241ng/ml

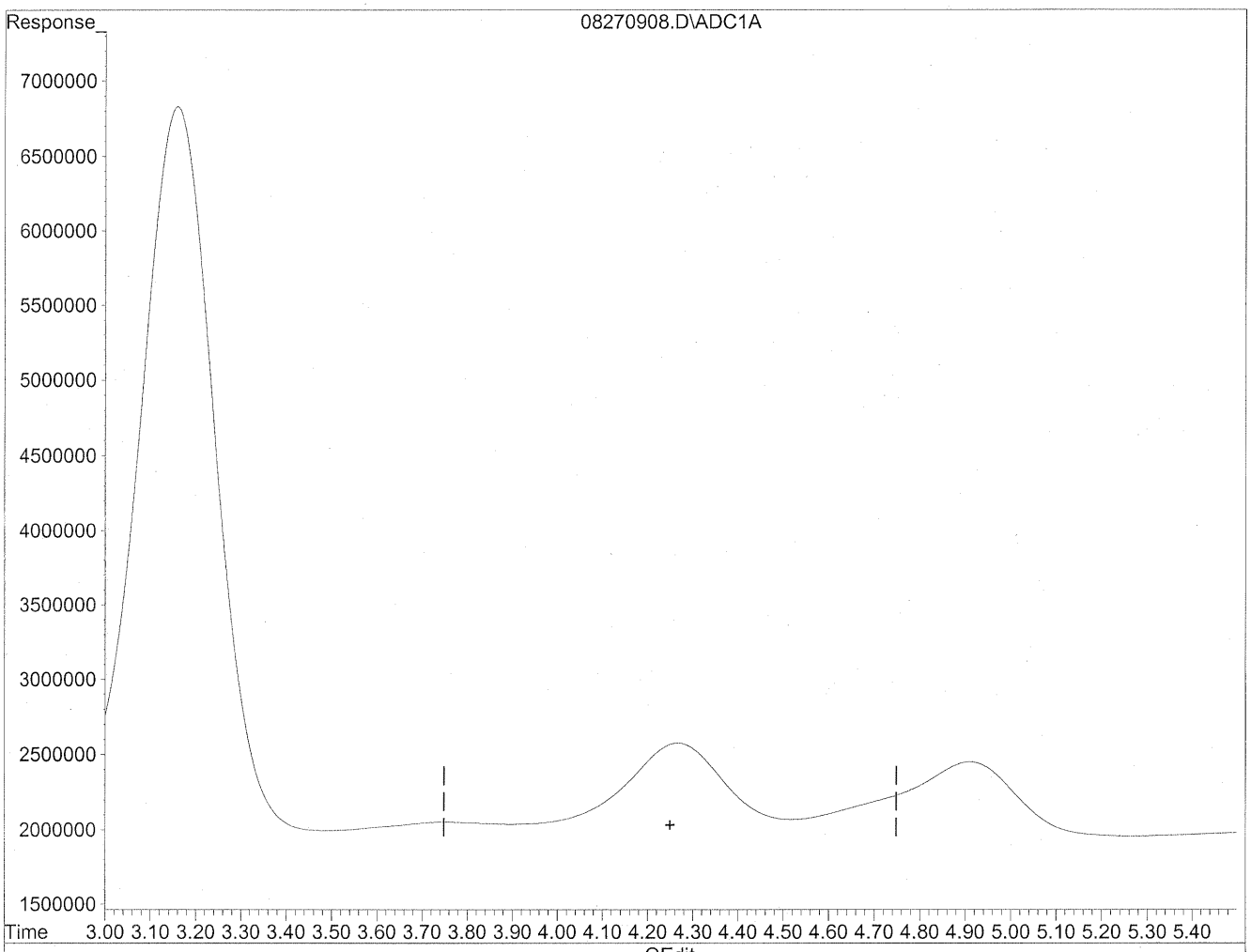
response 100750740

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



QEdit

(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/31/09 mm

LM 8/31/09

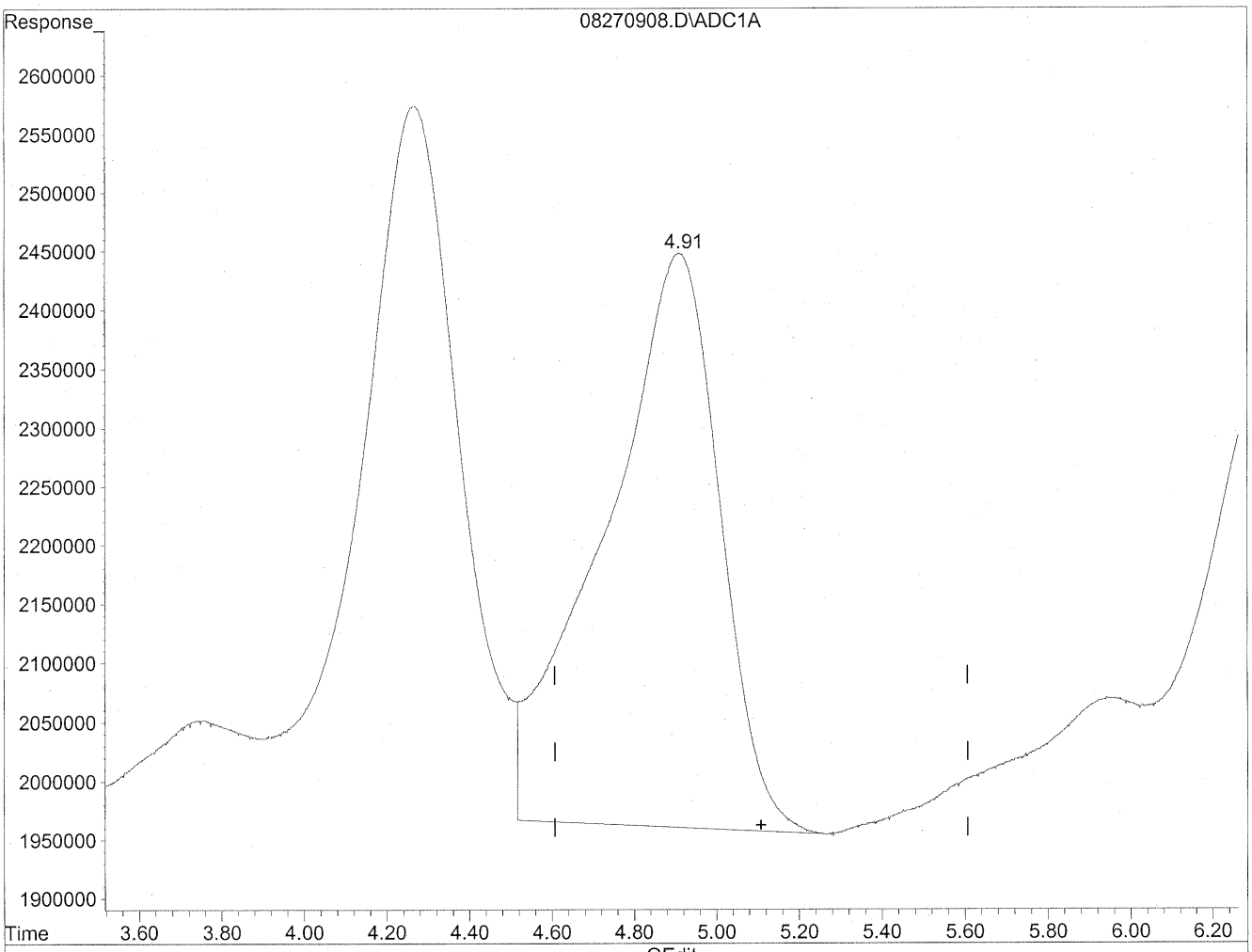
119

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.91min 1073.479ng/ml

response 94827010

120

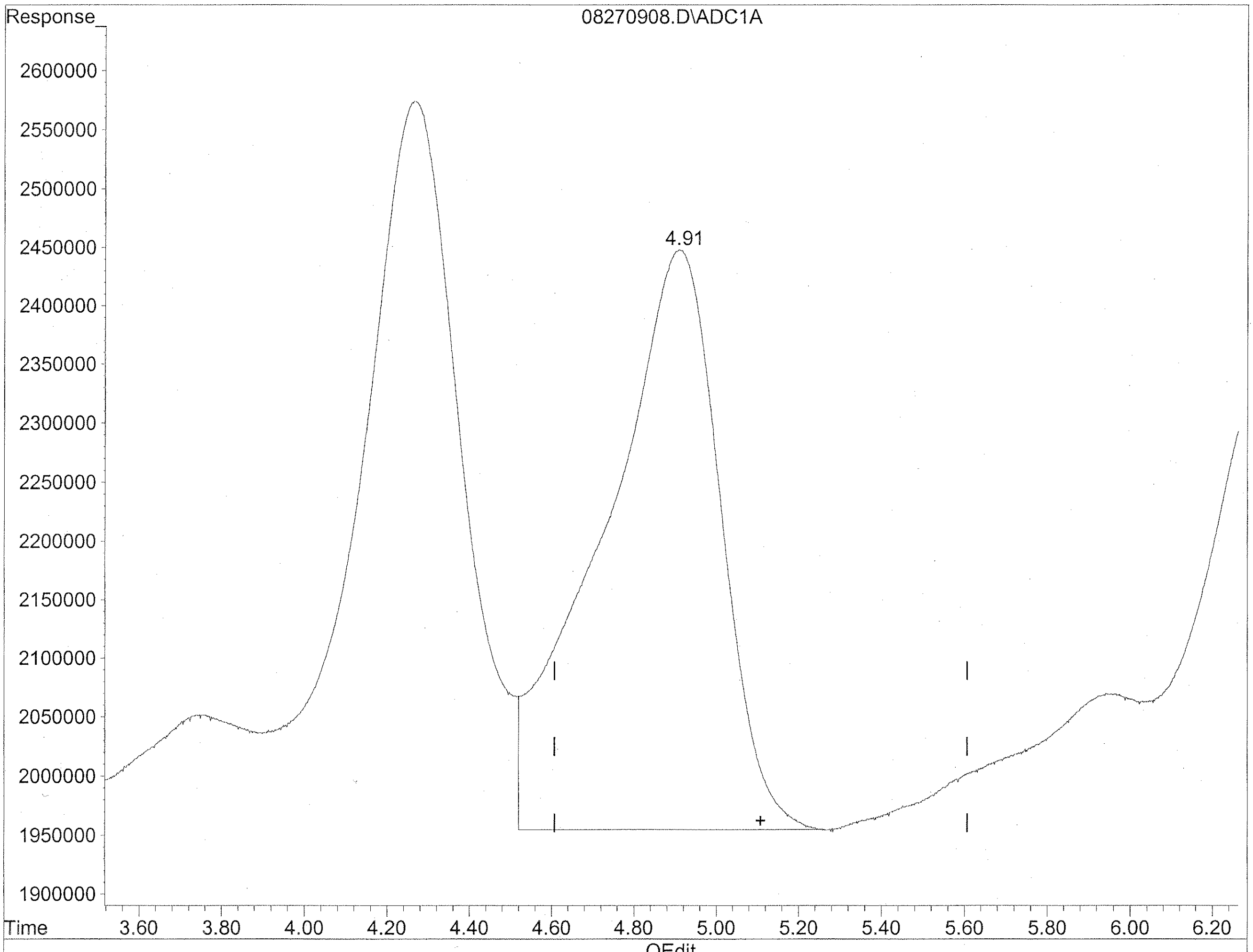
(+) = Expected Retention Time

08270908.D TO110709.M Sun Aug 30 13:37:35 2009

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D      Vial: 8
Acq On    : 27 Aug 2009  10:50 am                    Operator: HC
Sample    : P0902942-005 front 1.0ml                 Inst    : LC 01
Misc      :                                          Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Aug 27 11:38 19109  Quant Results File: TO110709.RES

Method        : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title         : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update   : Sat Aug 29 17:49:00 2009
Response via  : Multiple Level Calibration
```



(5) Butyraldehyde

4.91min 1100.646ng/ml m

response 97226838

HC 8/31/09 BC MP

LW 8/31/09

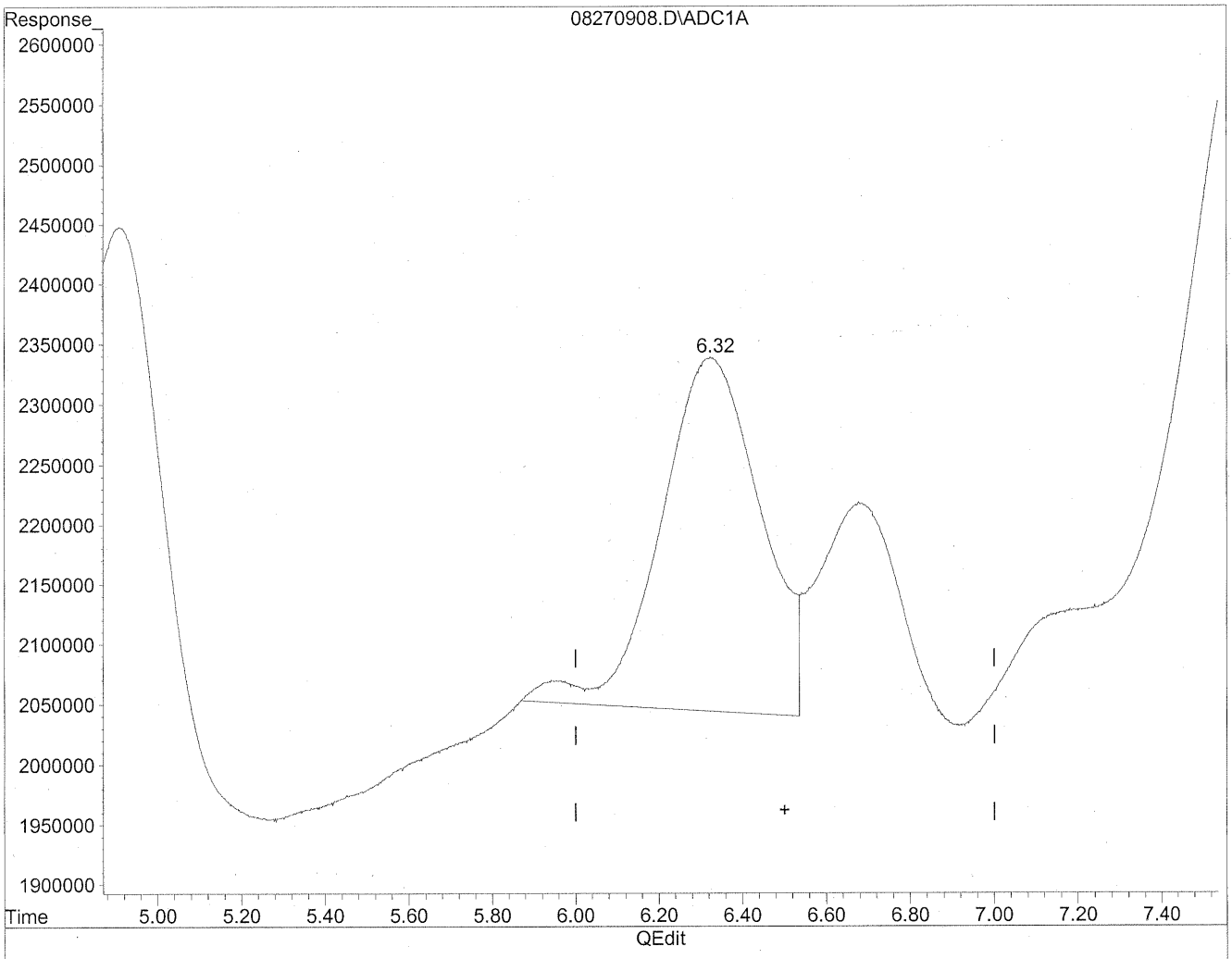
(+) = Expected Retention Time

08270908.D TO110709.M Sun Aug 30 13:37:50 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.32min 747.750ng/ml

response 49253750

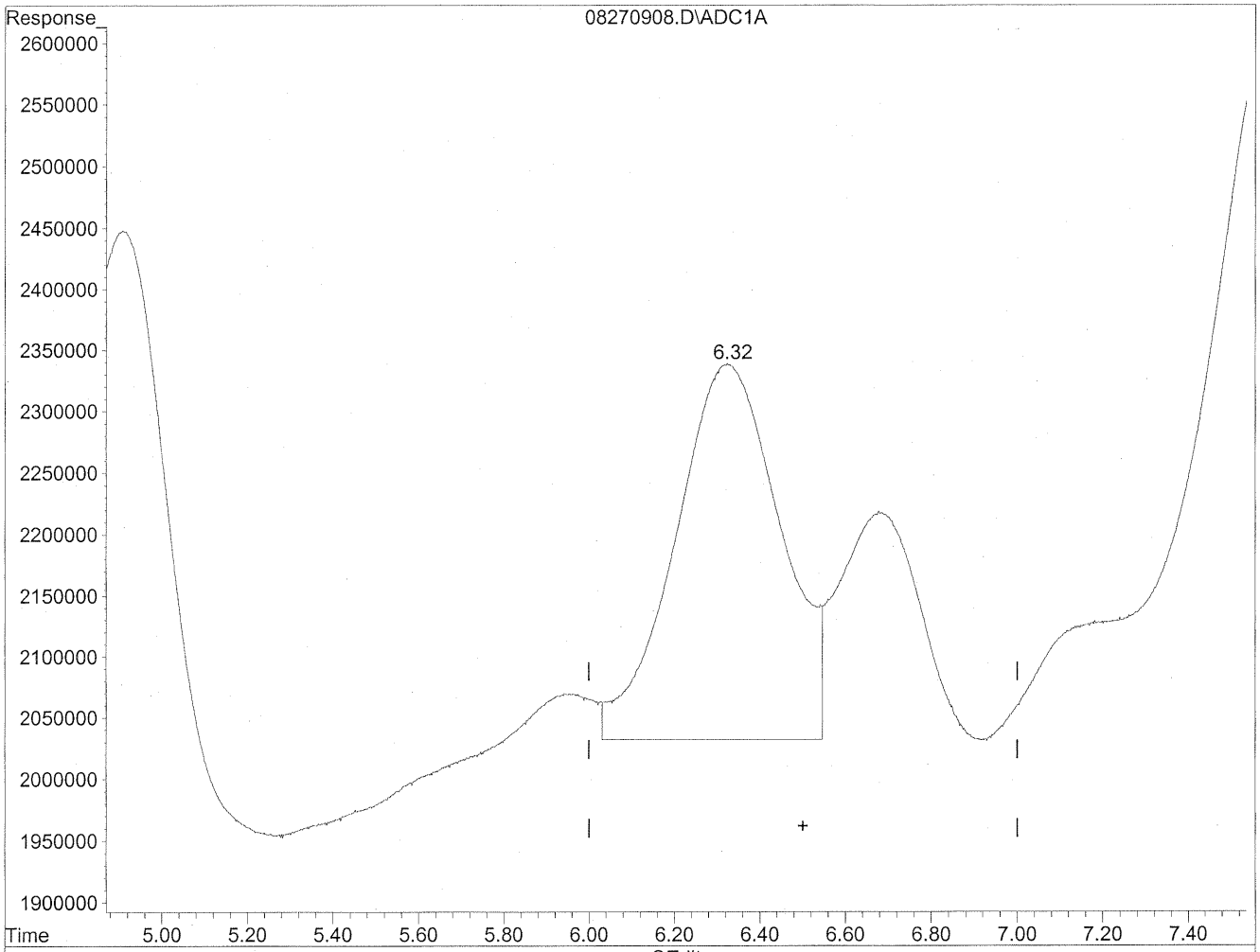
(+) = Expected Retention Time

08270908.D TO110709.M Sun Aug 30 13:37:54 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.32min 796.082ng/ml m

response 52437346

HC 8/31/09
BC

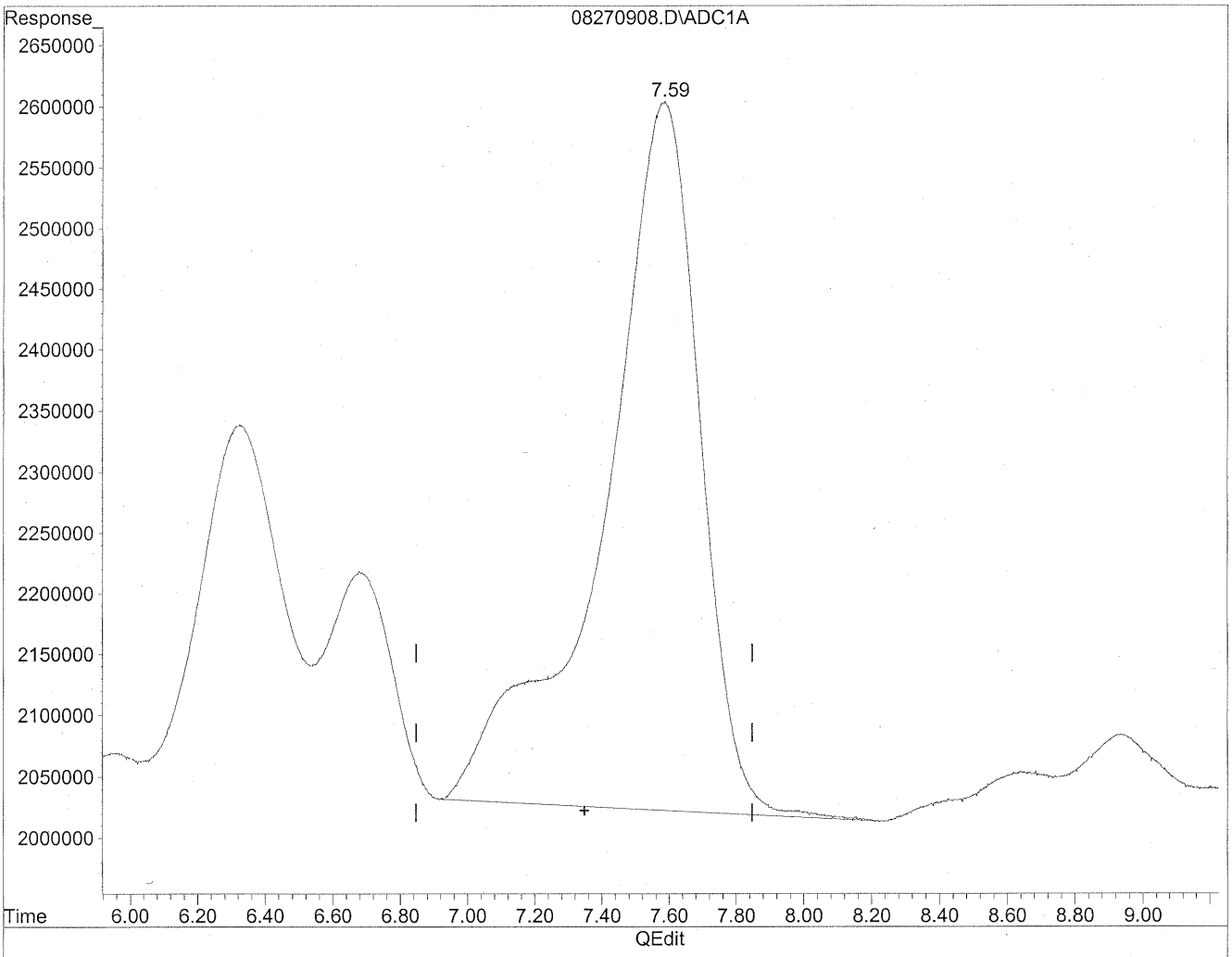
123

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.58min 1480.866ng/ml

response 115879261

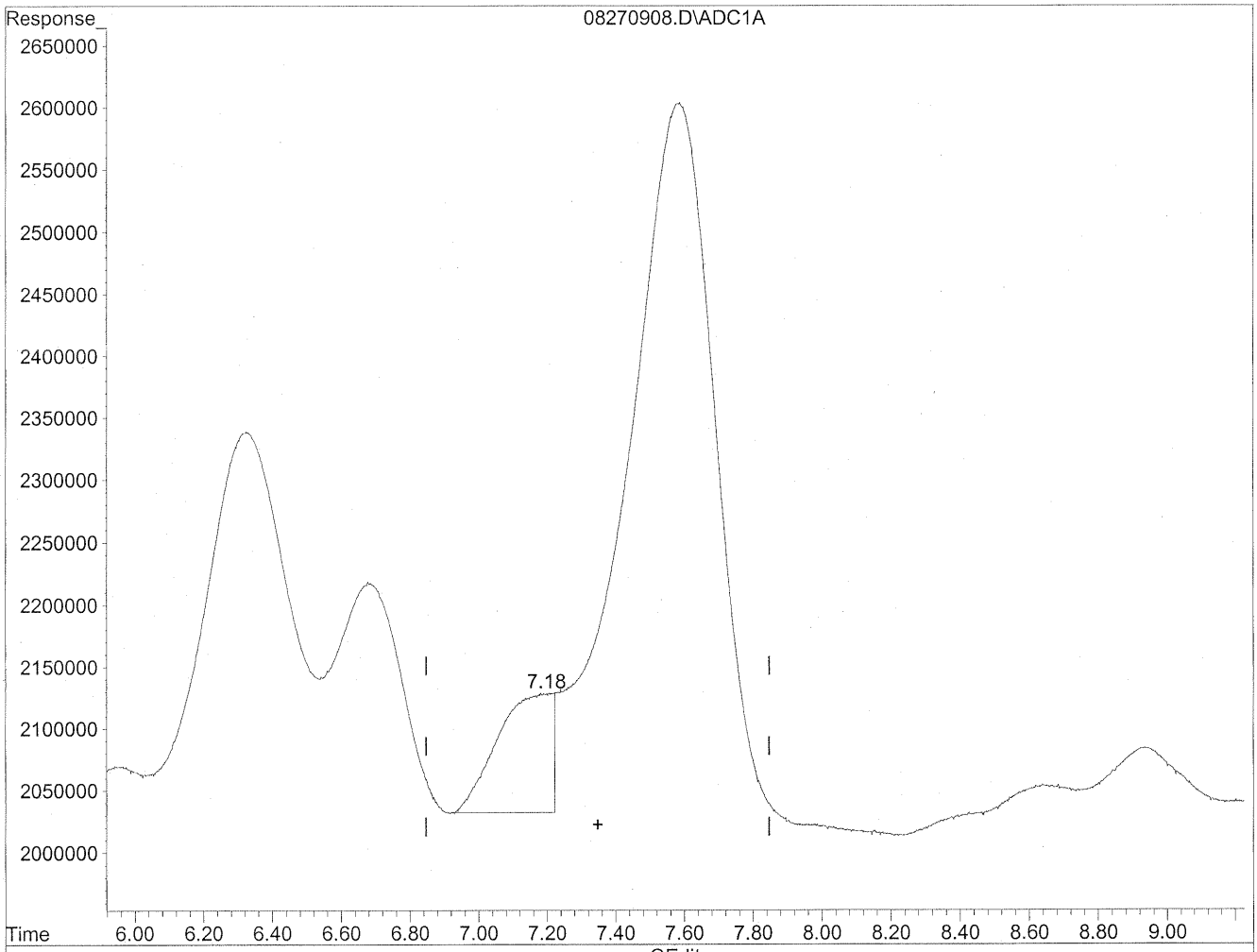
124

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.18min 137.045ng/ml m

response 10723919

HC 8/31/09
SH

8/31/09

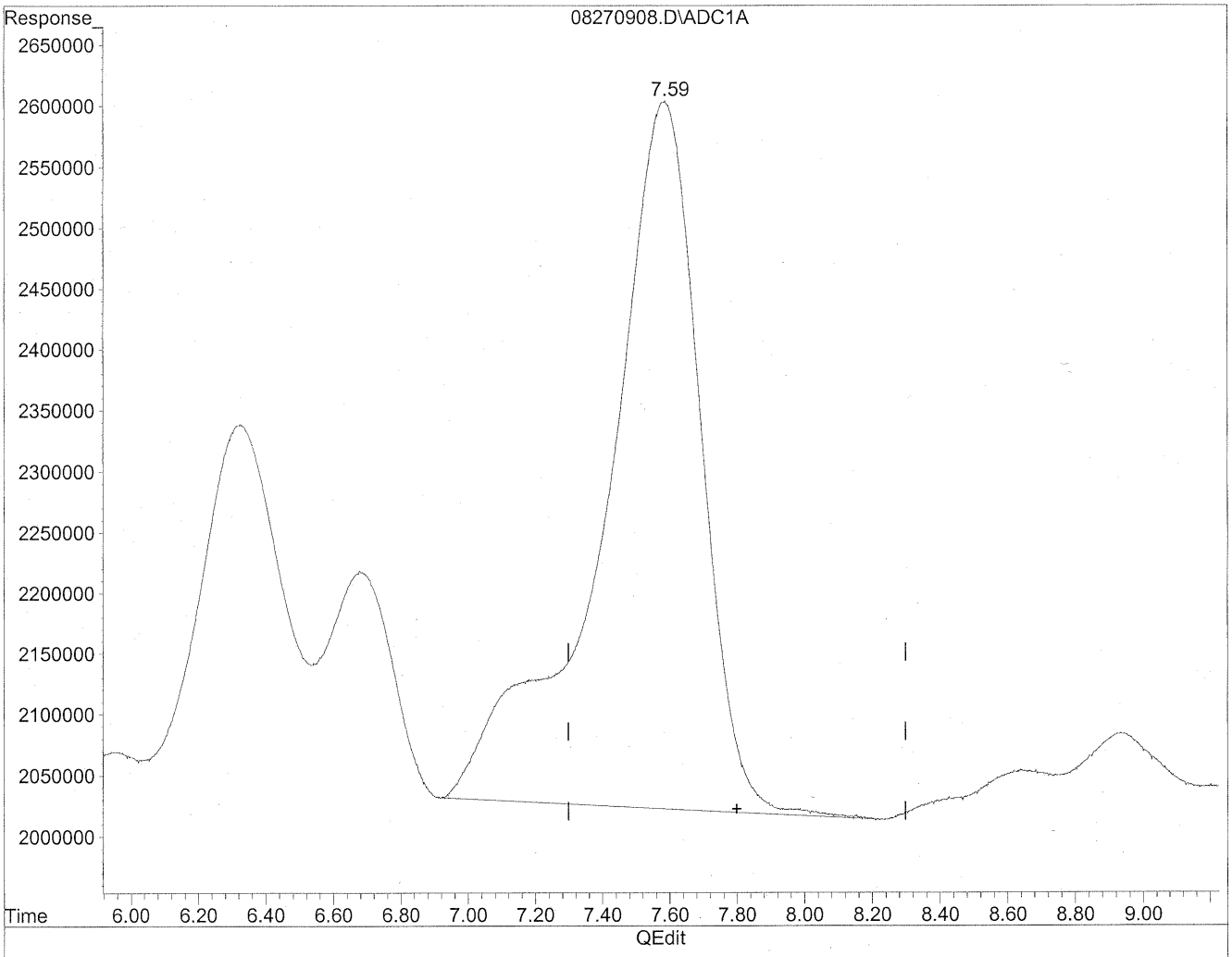
125

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.58min 1576.481ng/ml

response 115879261

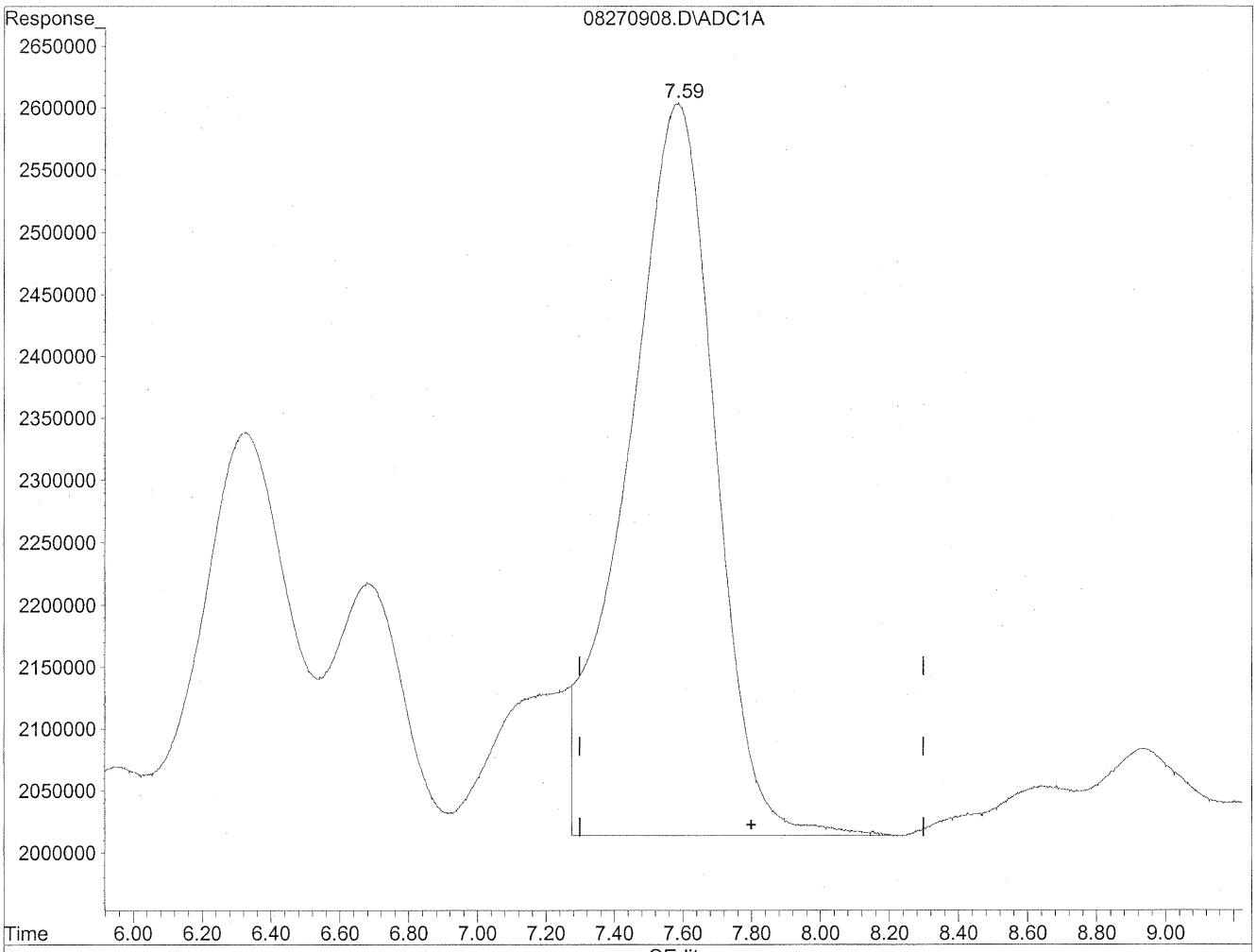
126

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.59min 1430.644ng/ml m

response 105159475

HC 8/31/09
SH/BC
WD 8/31/09

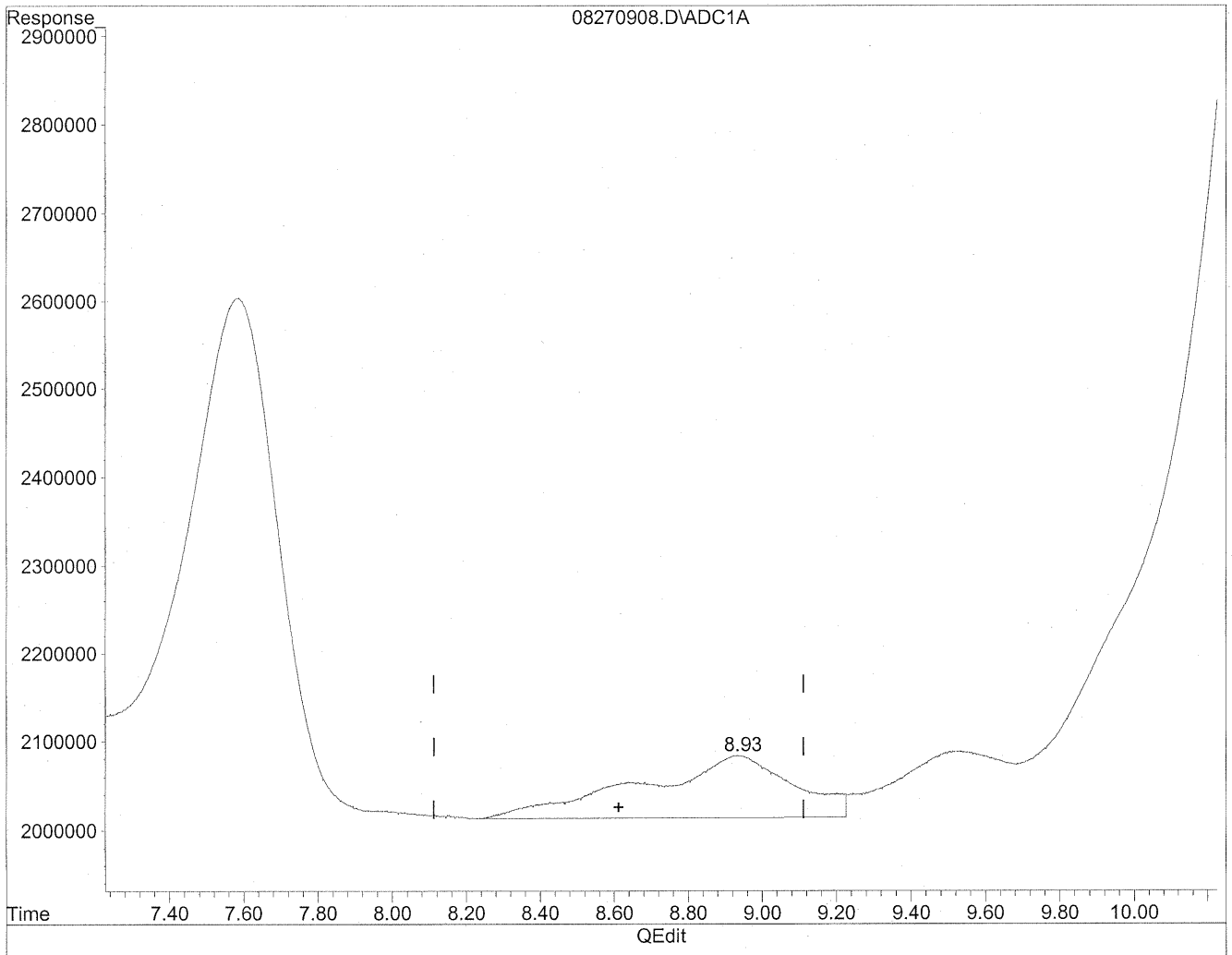
127

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

8.94min 340.683ng/ml

response 19868787

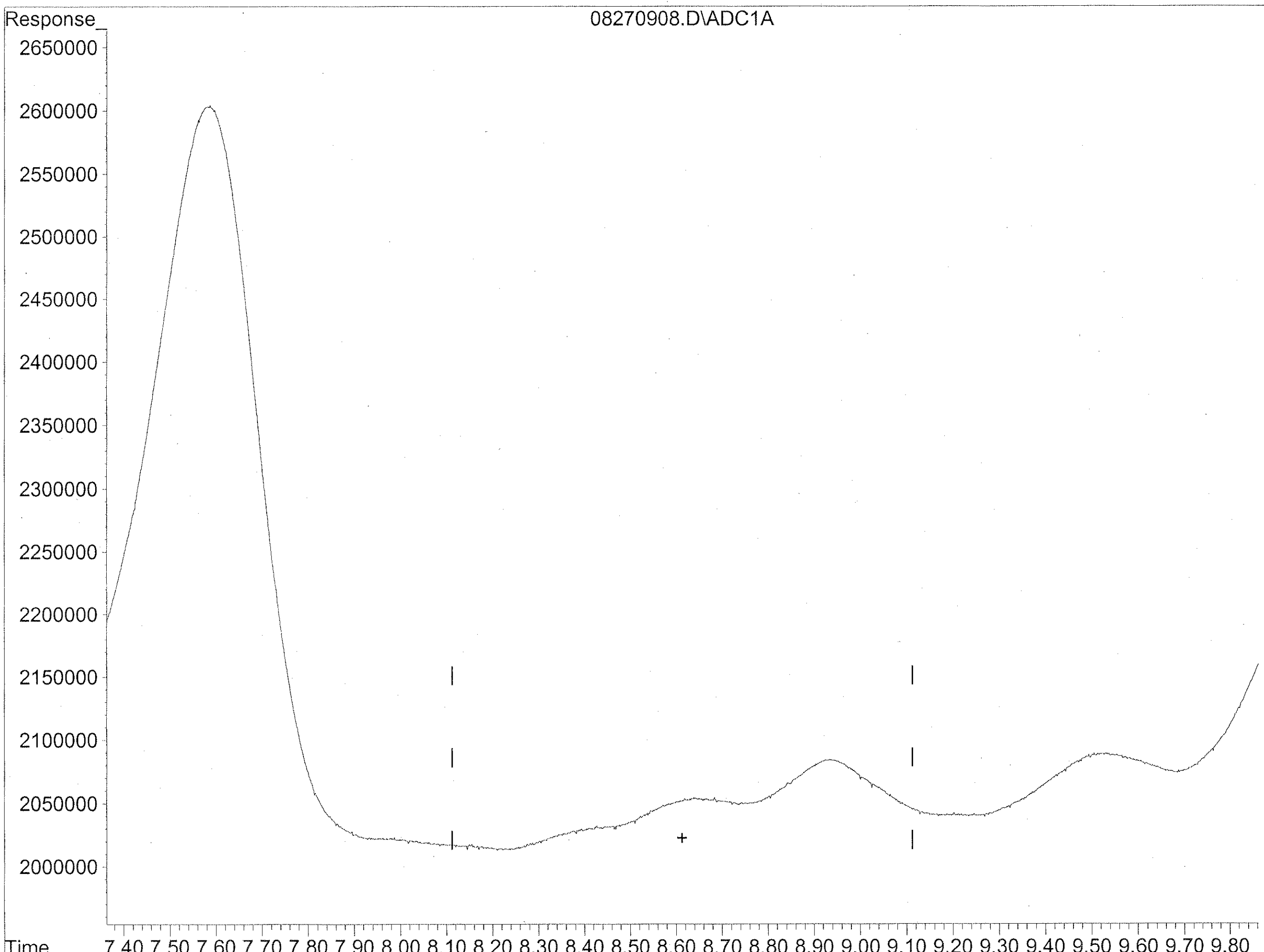
(+) = Expected Retention Time

08270908.D TO110709.M Sun Aug 30 13:38:41 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



QEdit

(9) o-Tolualdehyde

0.00min 0.000ng/ml d

response 0

HC 8/27/09 uup

ws 8/31/09

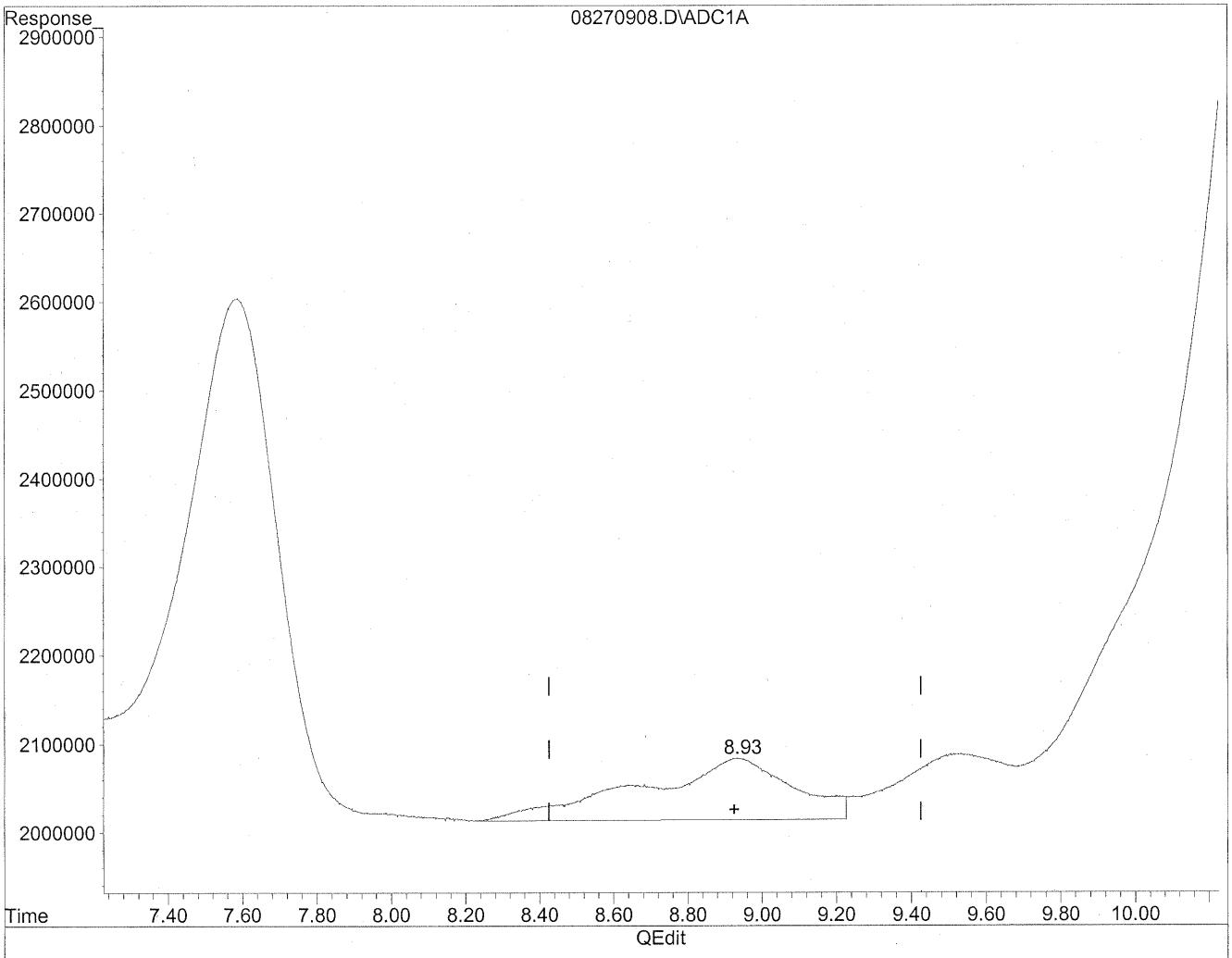
(+) = Expected Retention Time

08270908.D TO110709.M Sun Aug 30 13:38:44 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

8.94min 367.972ng/ml

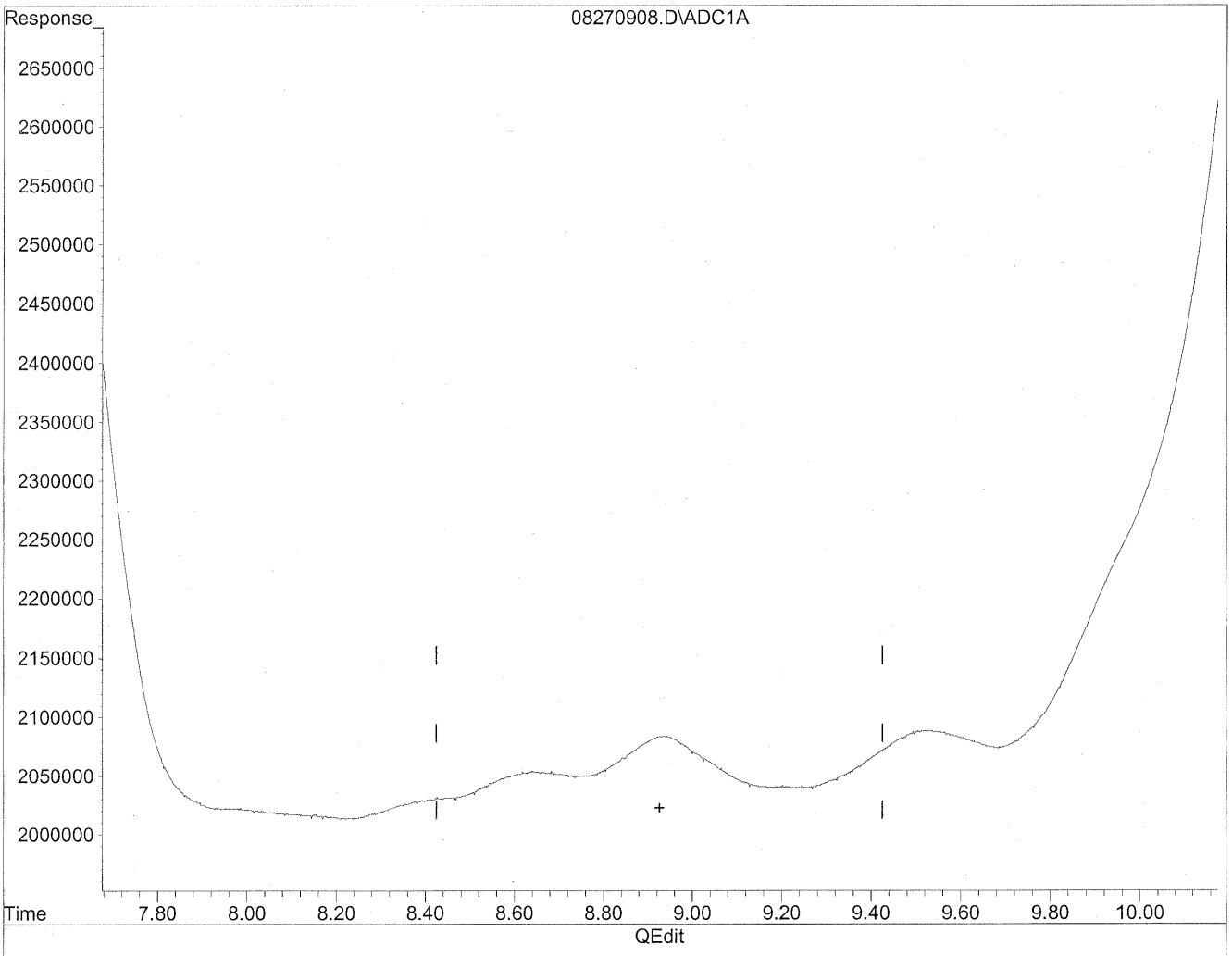
response 19868787

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

0.00min 0.000ng/ml d

response 0

tc 8/31/09 MP

Lot 8/31/09

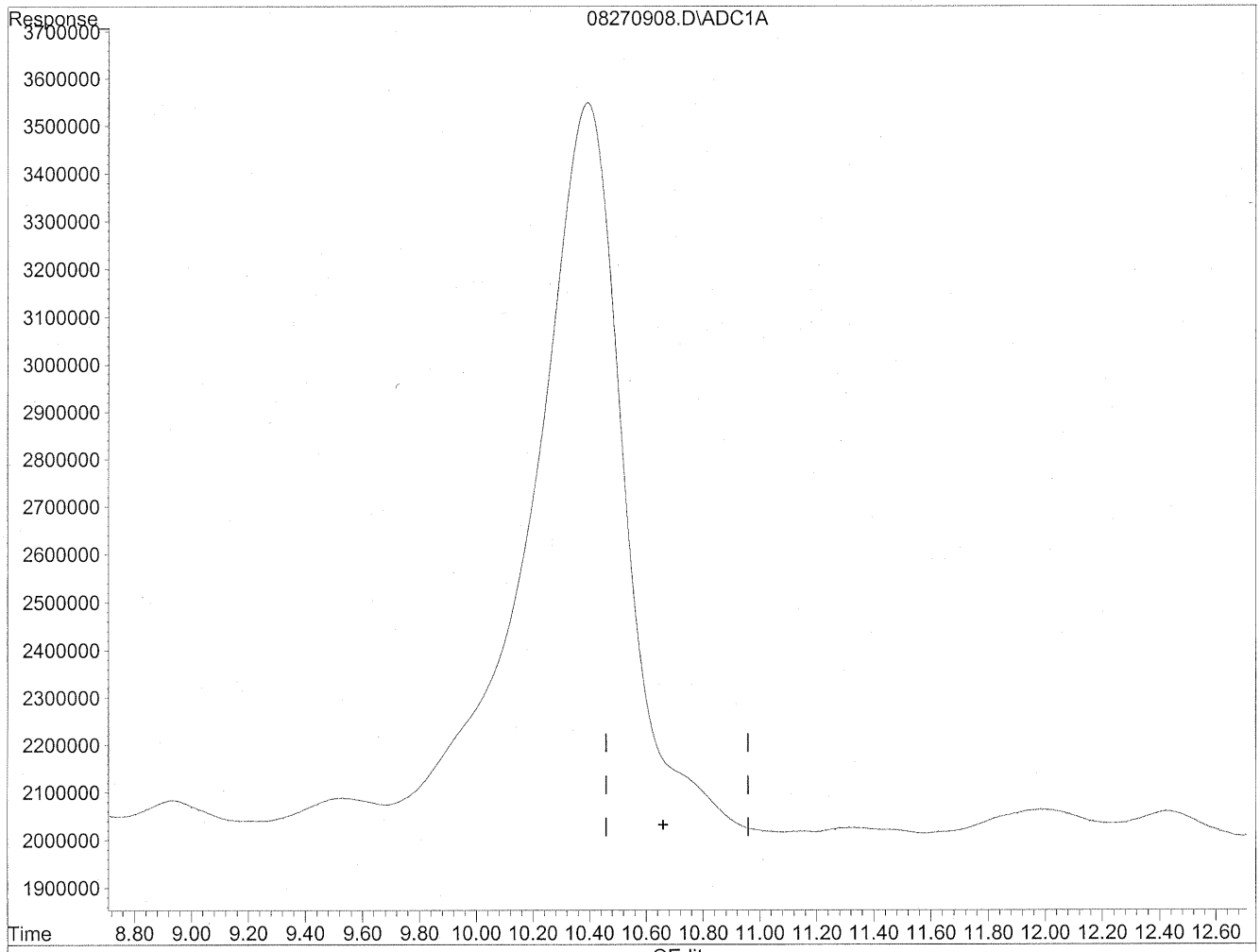
131

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



QEdit

(11) Hexaldehyde

10.66min 0.000ng/ml

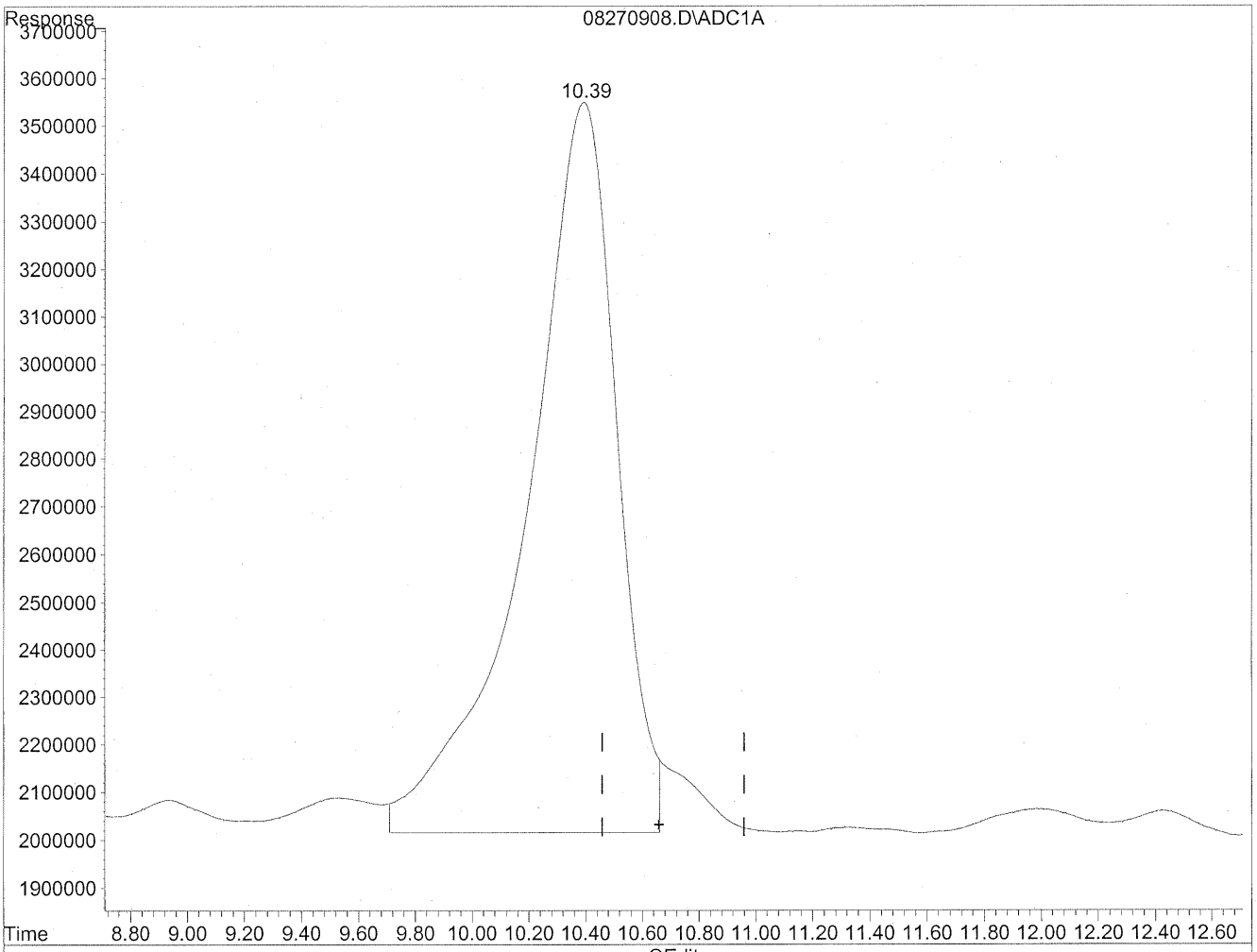
response 0

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270908.D Vial: 8
Acq On : 27 Aug 2009 10:50 am Operator: HC
Sample : P0902942-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.39min 5021.234ng/ml m

response 338148817

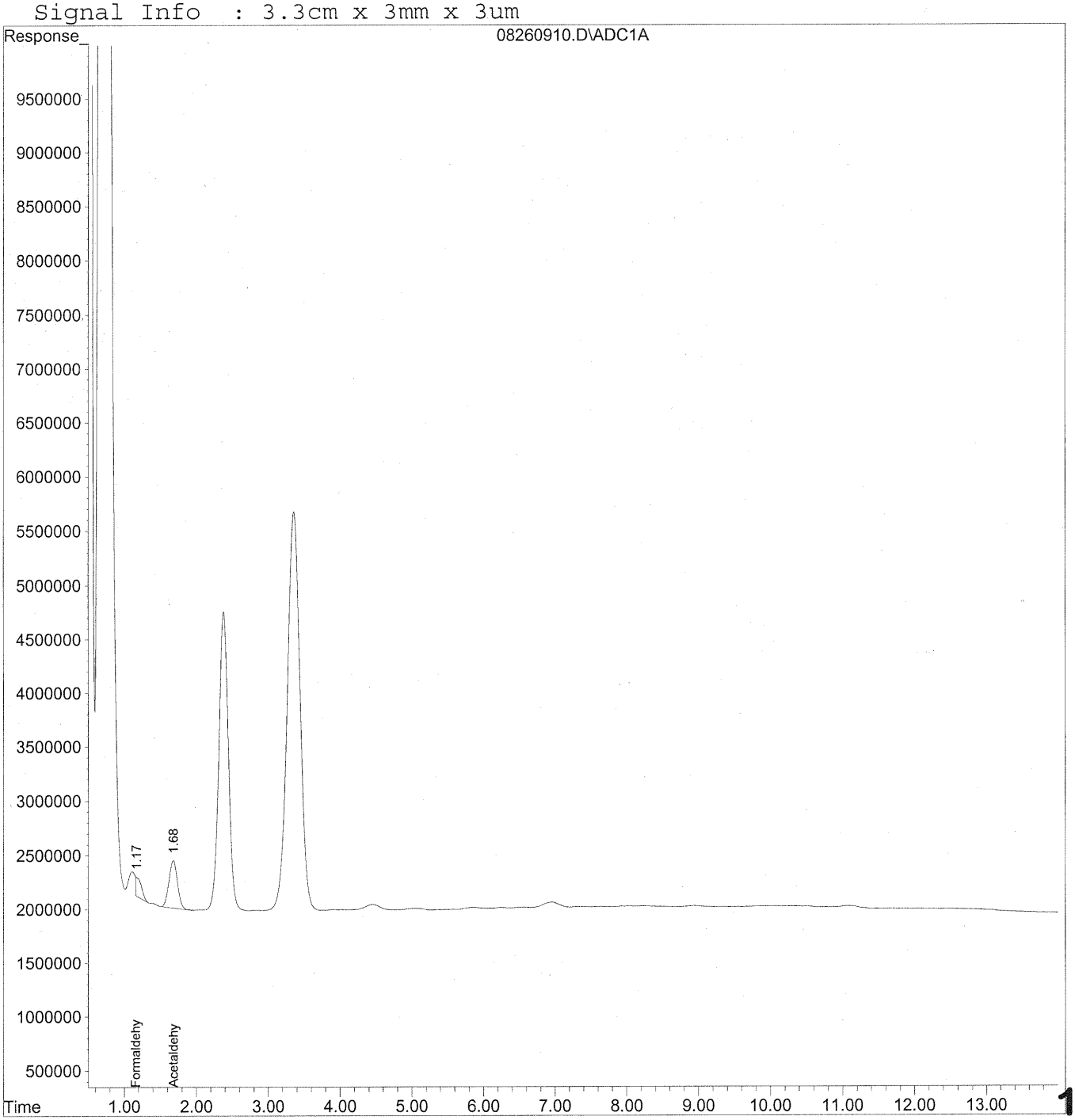
HC 8/31/09
BC 8/31/09

(+) = Expected Retention Time
08270908.D TO110709.M Sun Aug 30 13:39:29 2009

Data File : J:\LC01\DATA\TO11\2009_08\26\08260910.D Vial: 10
Acq On : 26 Aug 2009 7:20 pm Operator: HC
Sample : P0902942-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:50 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 : Thu Aug 27 07:35:56 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\26\08260910.D Vial: 10
Acq On : 26 Aug 2009 7:20 pm Operator: HC
Sample : P0902942-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:50 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 : Thu Aug 27 07:35:56 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

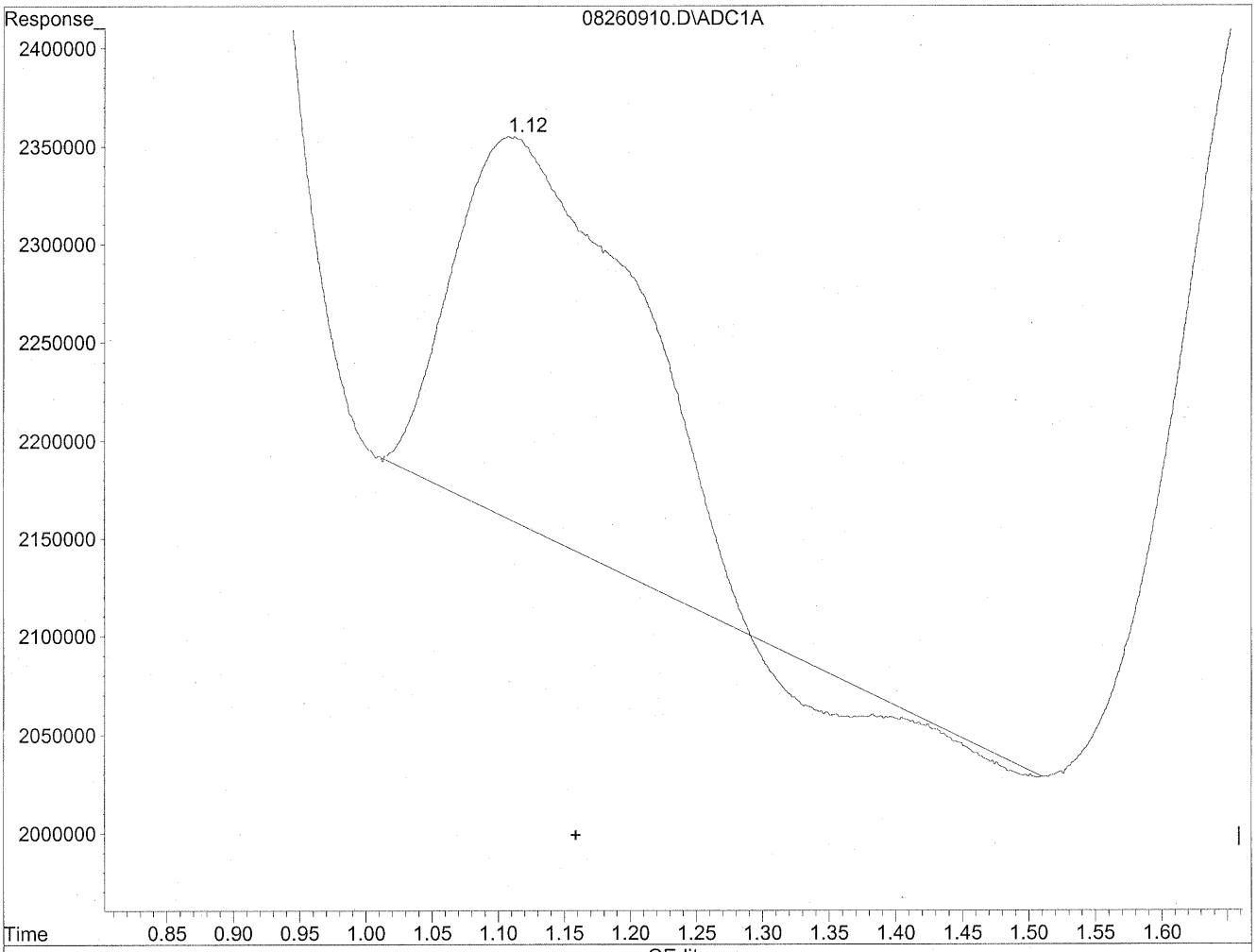
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.17	9628255	52.447	ng/mlm
2)	Acetaldehyde	1.68	36422867	259.749	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\26\08260910.D Vial: 10
Acq On : 26 Aug 2009 7:20 pm Operator: HC
Sample : P0902942-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.11min 100.501ng/ml

response 18450170

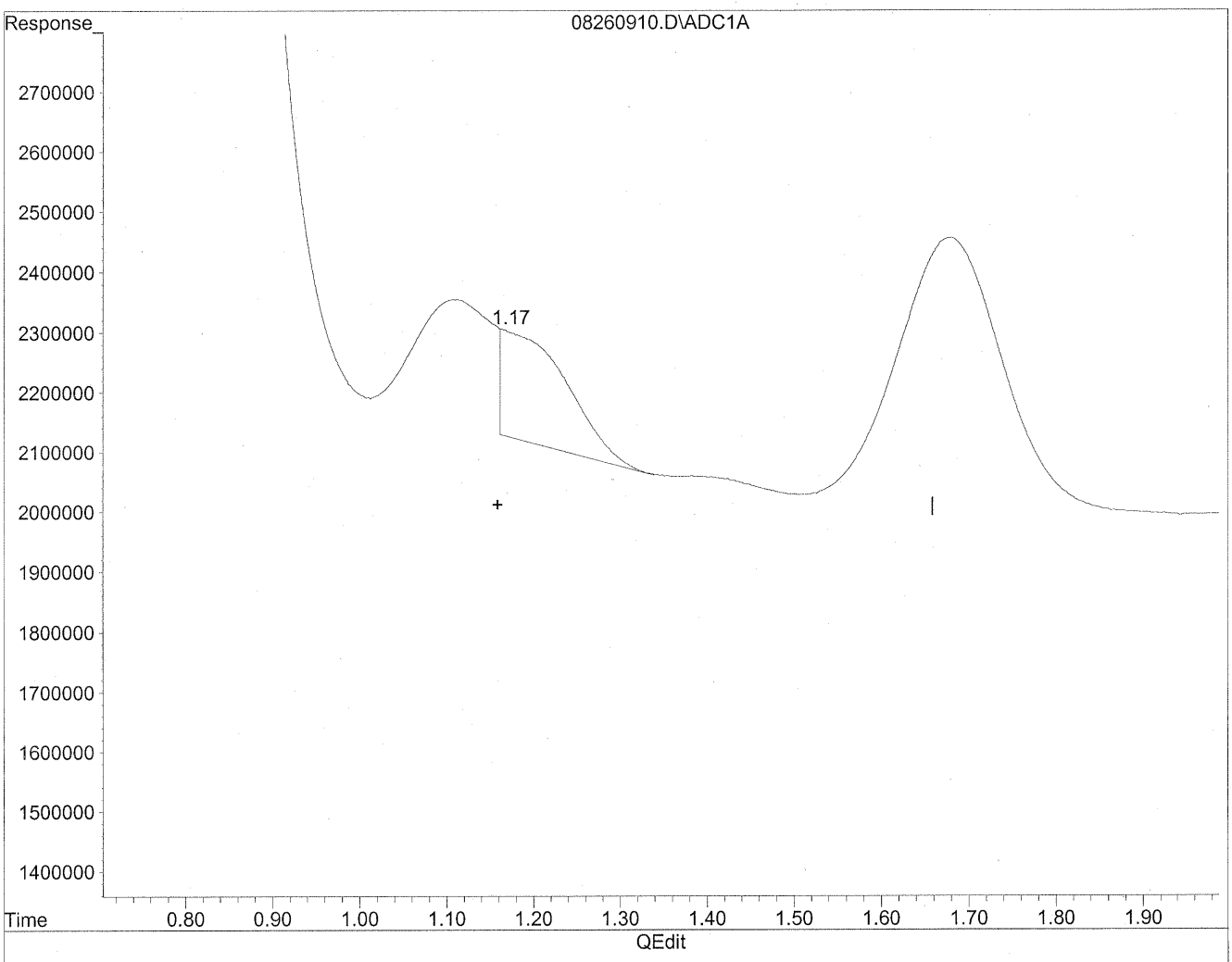
(+) = Expected Retention Time

08260910.D TO110709.M Sat Aug 29 16:50:04 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260910.D Vial: 10
Acq On : 26 Aug 2009 7:20 pm Operator: HC
Sample : P0902942-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.17min 52.447ng/ml m

response 9628255

HC 8/30/09

SP

8/31/09

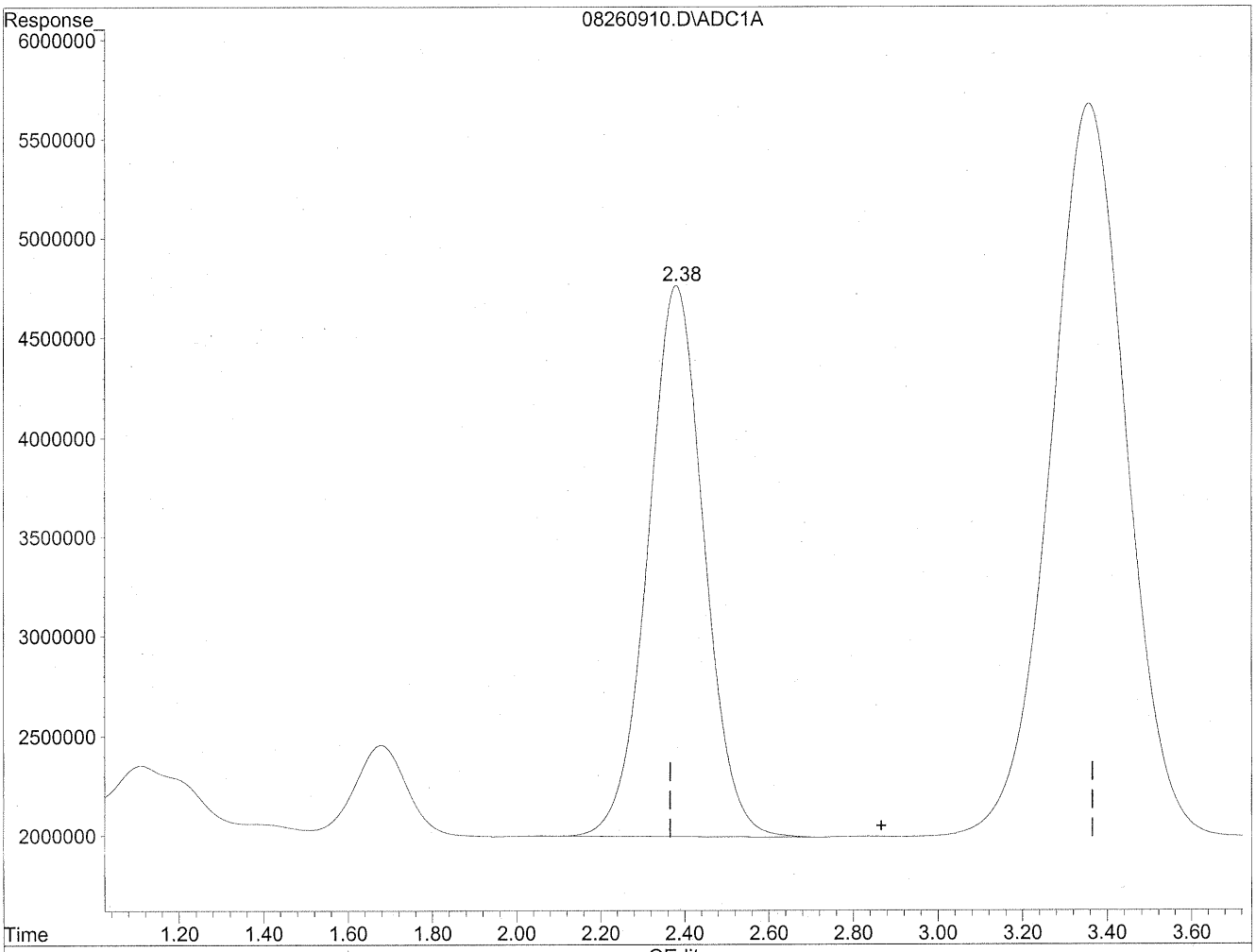
(+) = Expected Retention Time

08260910.D TO110709.M Sat Aug 29 16:50:29 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260910.D Vial: 10
Acq On : 26 Aug 2009 7:20 pm Operator: HC
Sample : P0902942-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.38min 2404.624ng/ml

response 256561897

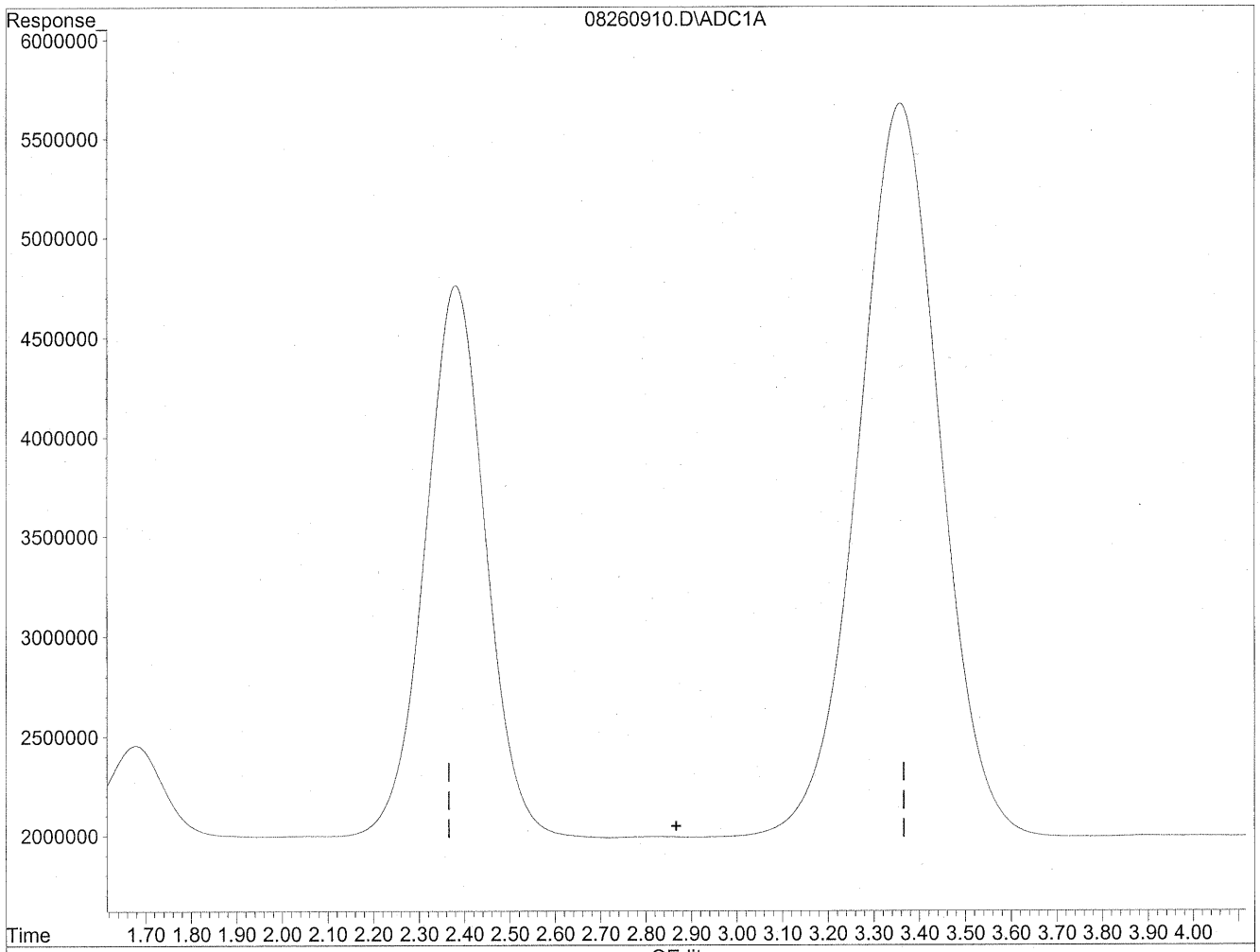
138

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260910.D Vial: 10
Acq On : 26 Aug 2009 7:20 pm Operator: HC
Sample : P0902942-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



QEdit

(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/30/09 WP

WF 8/31/09

139

(+) = Expected Retention Time

08260910.D TO110709.M Sat Aug 29 16:50:45 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **101474**
Client Project ID: **16512**

CAS Project ID: P0902942
CAS Sample ID: P0902942-006

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: 8/24/09
Date Received: 8/25/09
Date Analyzed: 8/26-27/09
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	< 200	NA	NA	NA	NA	V
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	< 200	NA	NA	NA	NA	V

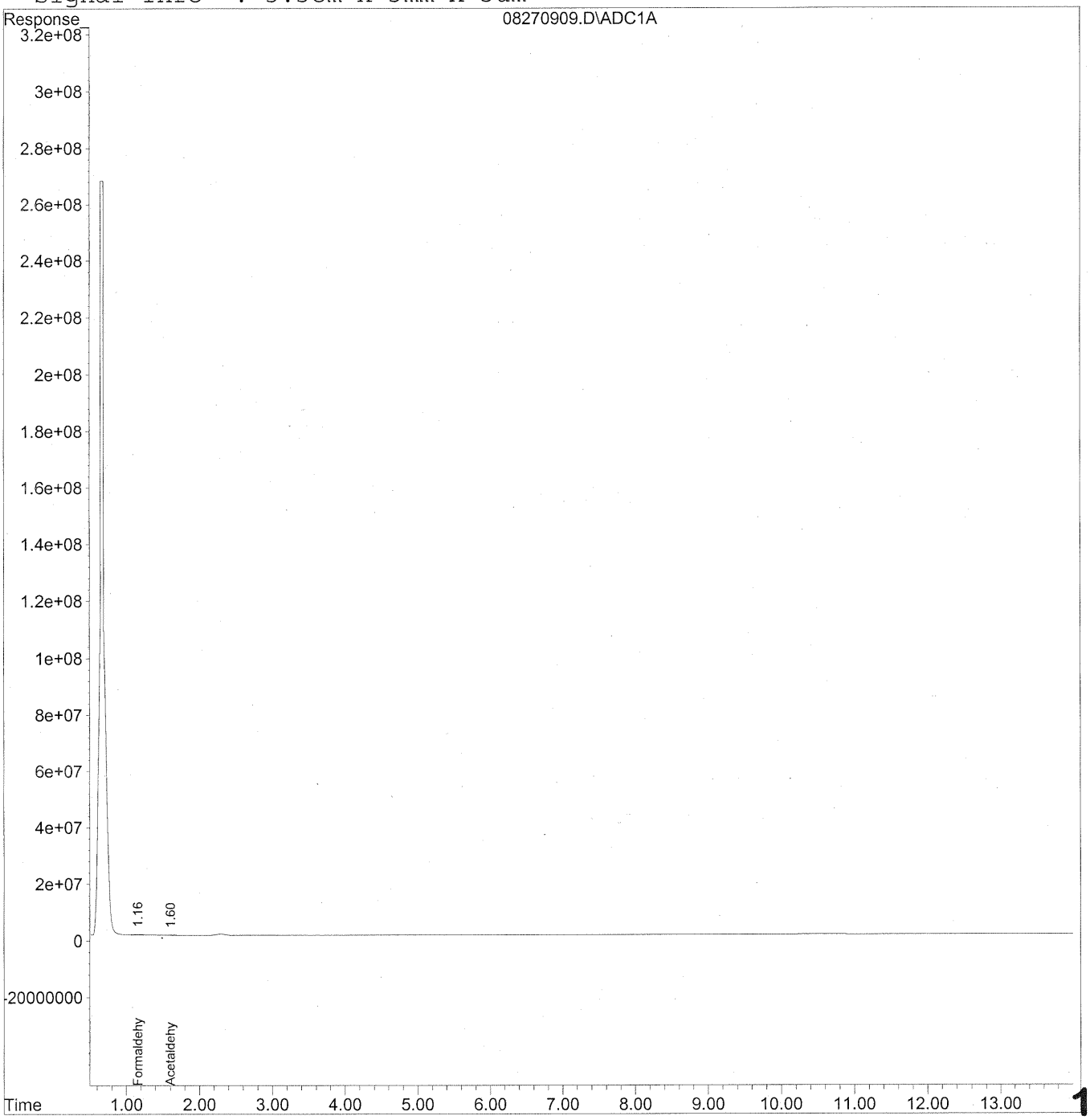
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
NA = Not applicable.
V = The continuing calibration verification standard was outside (biased low) the specified limits for this compound.

Verified By: Rc. Date: 9/11/09

Data File : J:\LC01\DATA\TO11\2009_08\27\08270909.D Vial: 9
Acq On : 27 Aug 2009 11:06 am Operator: HC
Sample : P0902942-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:40 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 : Thu Aug 27 08:33:51 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



141

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\27\08270909.D Vial: 9
Acq On : 27 Aug 2009 11:06 am Operator: HC
Sample : P0902942-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 30 13:40 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 : Thu Aug 27 08:33:51 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

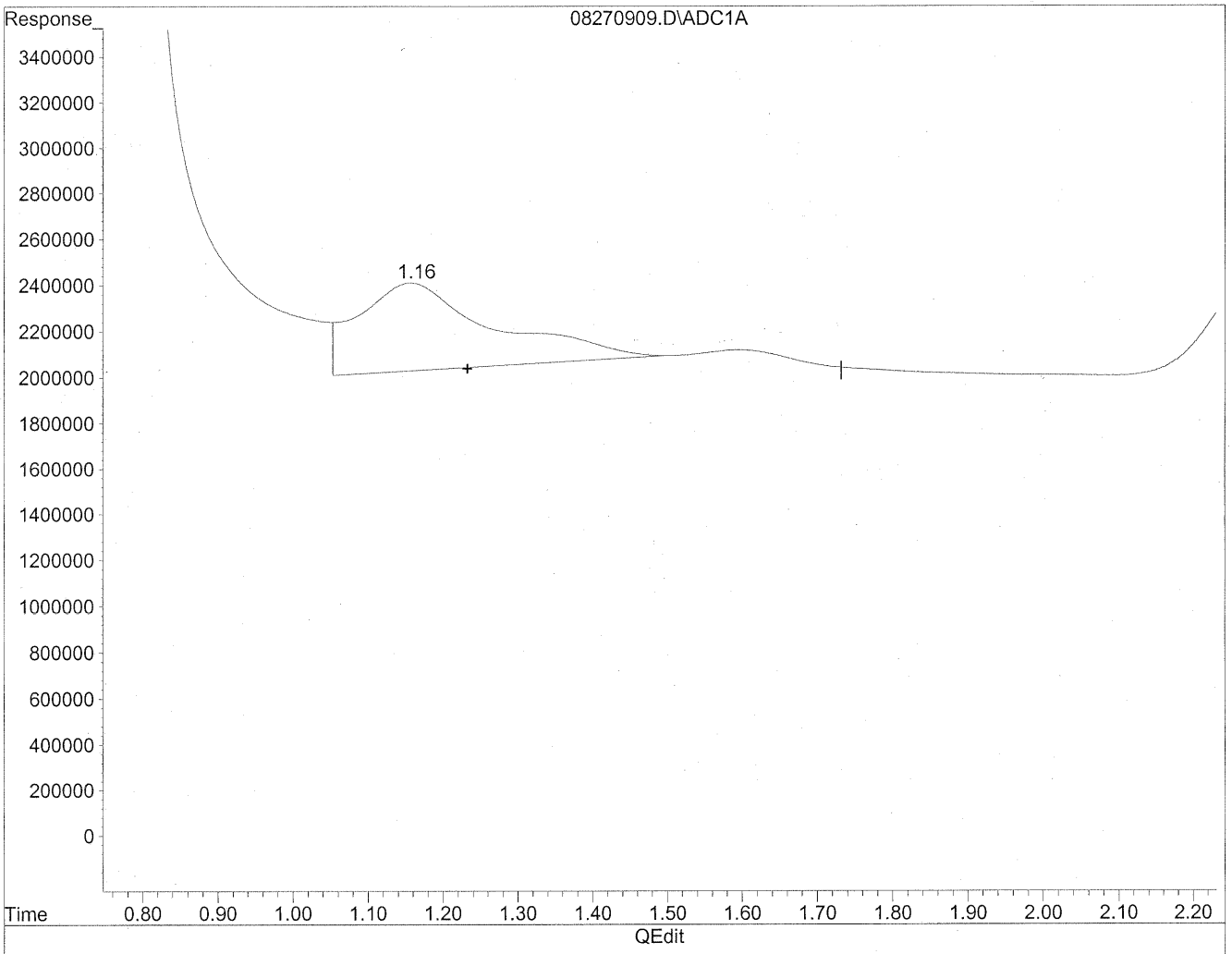
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	12230495	66.622	ng/mlm
2)	Acetaldehyde	1.60	2845731	20.294	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/mld
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270909.D Vial: 9
Acq On : 27 Aug 2009 11:06 am Operator: HC
Sample : P0902942-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.16min 258.885ng/ml

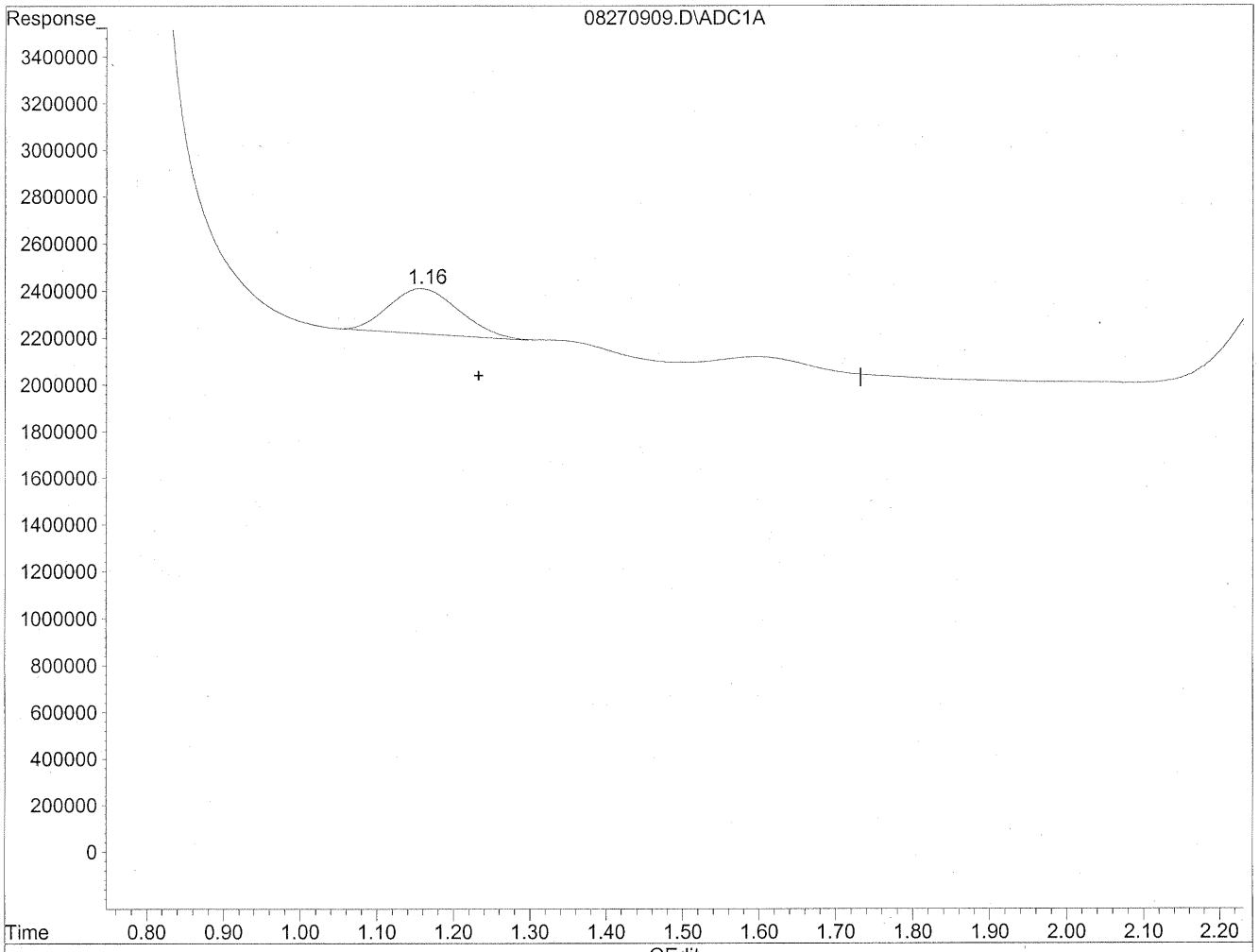
response 47526399

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270909.D Vial: 9
Acq On : 27 Aug 2009 11:06 am Operator: HC
Sample : P0902942-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.16min 66.622ng/ml m

response 12230495

HC 8/31/09 VL

8/31/09

144

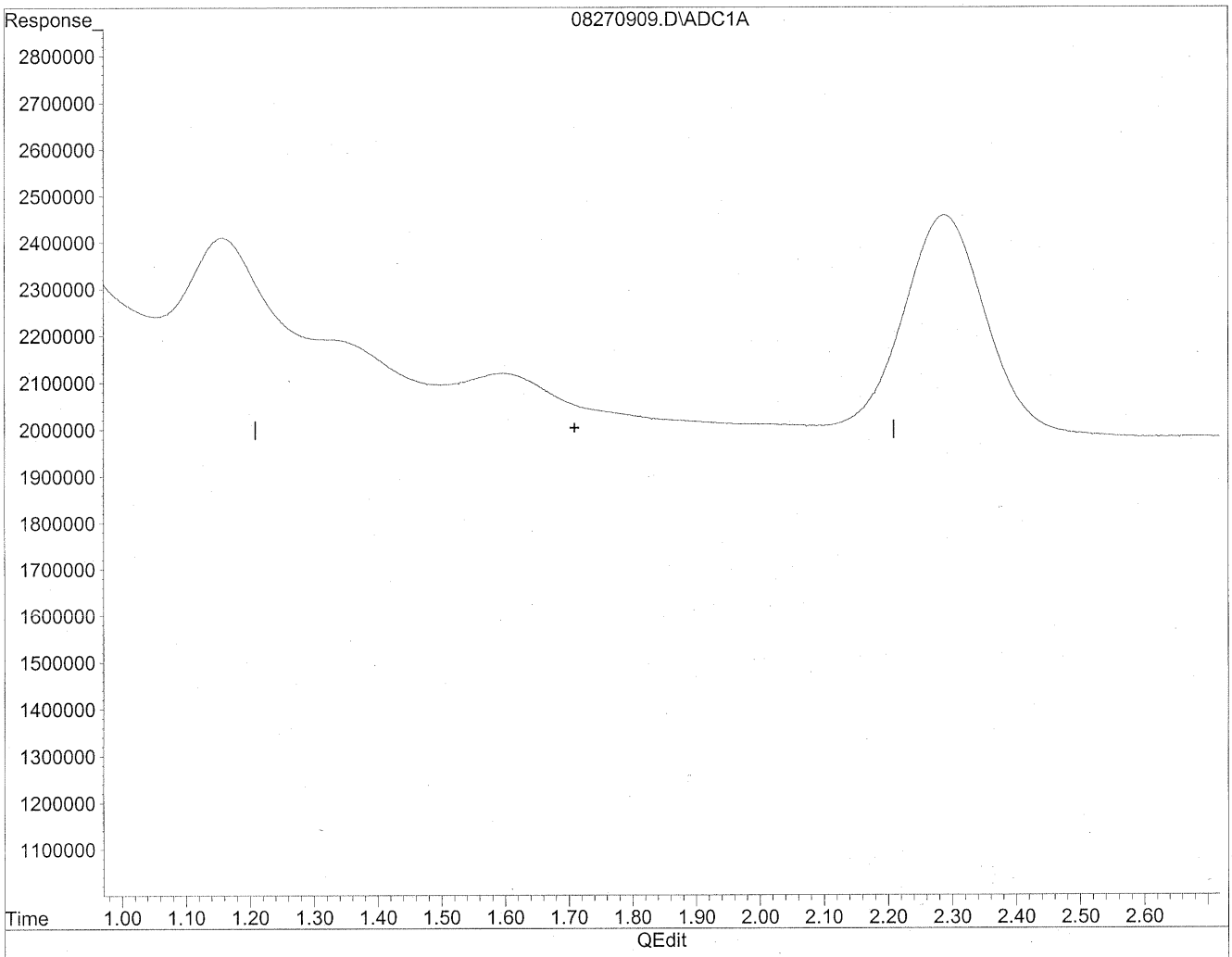
(+) = Expected Retention Time

08270909.D TO110709.M Sun Aug 30 13:39:46 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270909.D Vial: 9
Acq On : 27 Aug 2009 11:06 am Operator: HC
Sample : P0902942-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.71min 0.000ng/ml

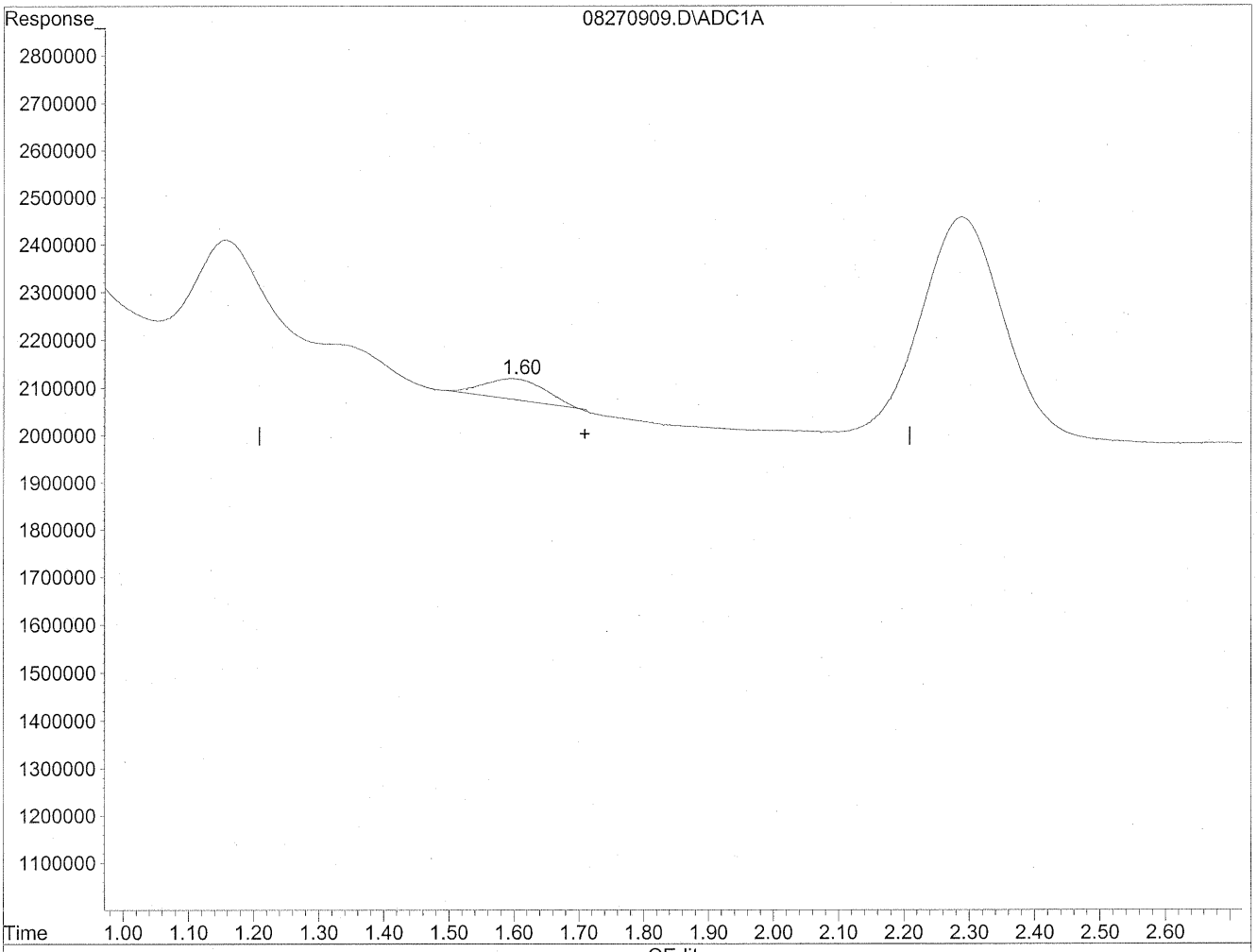
response 0

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270909.D Vial: 9
Acq On : 27 Aug 2009 11:06 am Operator: HC
Sample : P0902942-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



QEdit

(2) Acetaldehyde

1.60min 20.294ng/ml m

response 2845731

HL 8/31/09 LC

Log 8/31/09

146

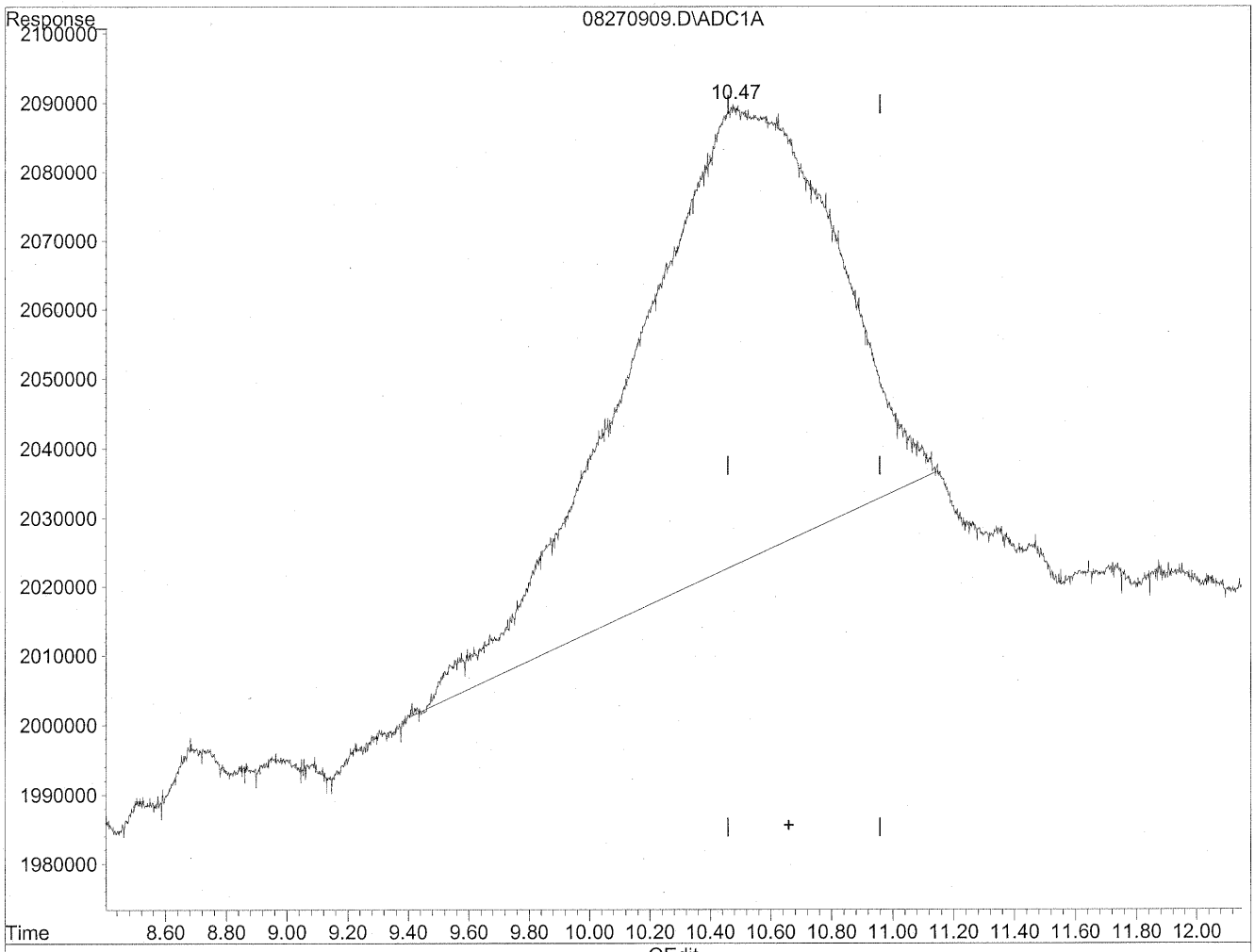
(+) = Expected Retention Time

08270909.D TO110709.M Sun Aug 30 13:40:00 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270909.D Vial: 9
Acq On : 27 Aug 2009 11:06 am Operator: HC
Sample : P0902942-006 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 11:38 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.48min 459.380ng/ml

response 30936410

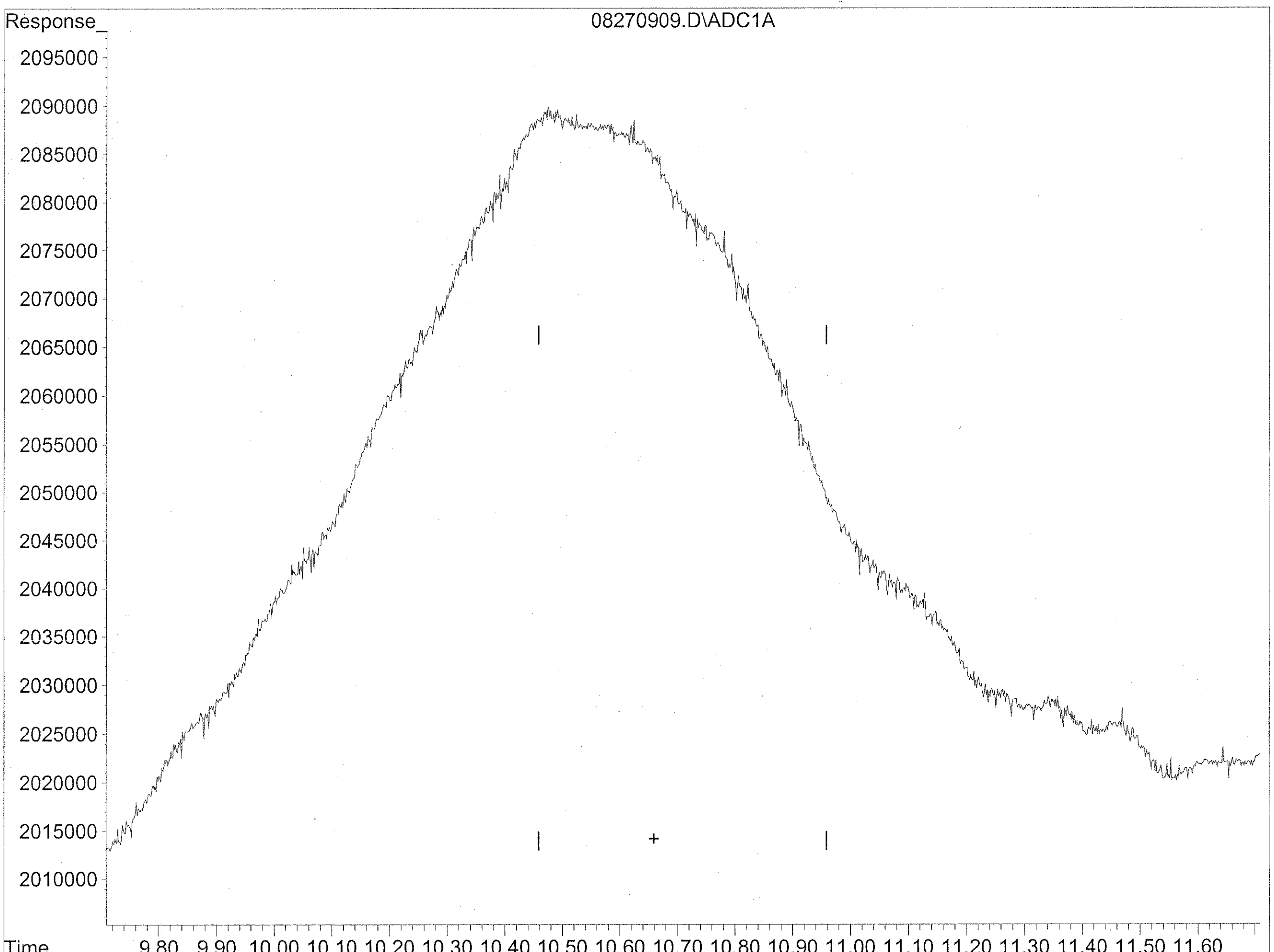
(+) = Expected Retention Time

08270909.D TO110709.M Sun Aug 30 13:40:07 2009

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2009_08\27\08270909.D       Vial: 9
Acq On    : 27 Aug 2009  11:06 am                      Operator: HC
Sample    : P0902942-006 front 1.0ml                   Inst    : LC 01
Misc      :                                            Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Aug 27 11:38 19109  Quant Results File: TO110709.RES

Method        : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title         : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update   : Sat Aug 29 17:49:00 2009
Response via  : Multiple Level Calibration
```



QEdit

(11) Hexaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/31/09 no peak

VH 8/31/09

(+) = Expected Retention Time

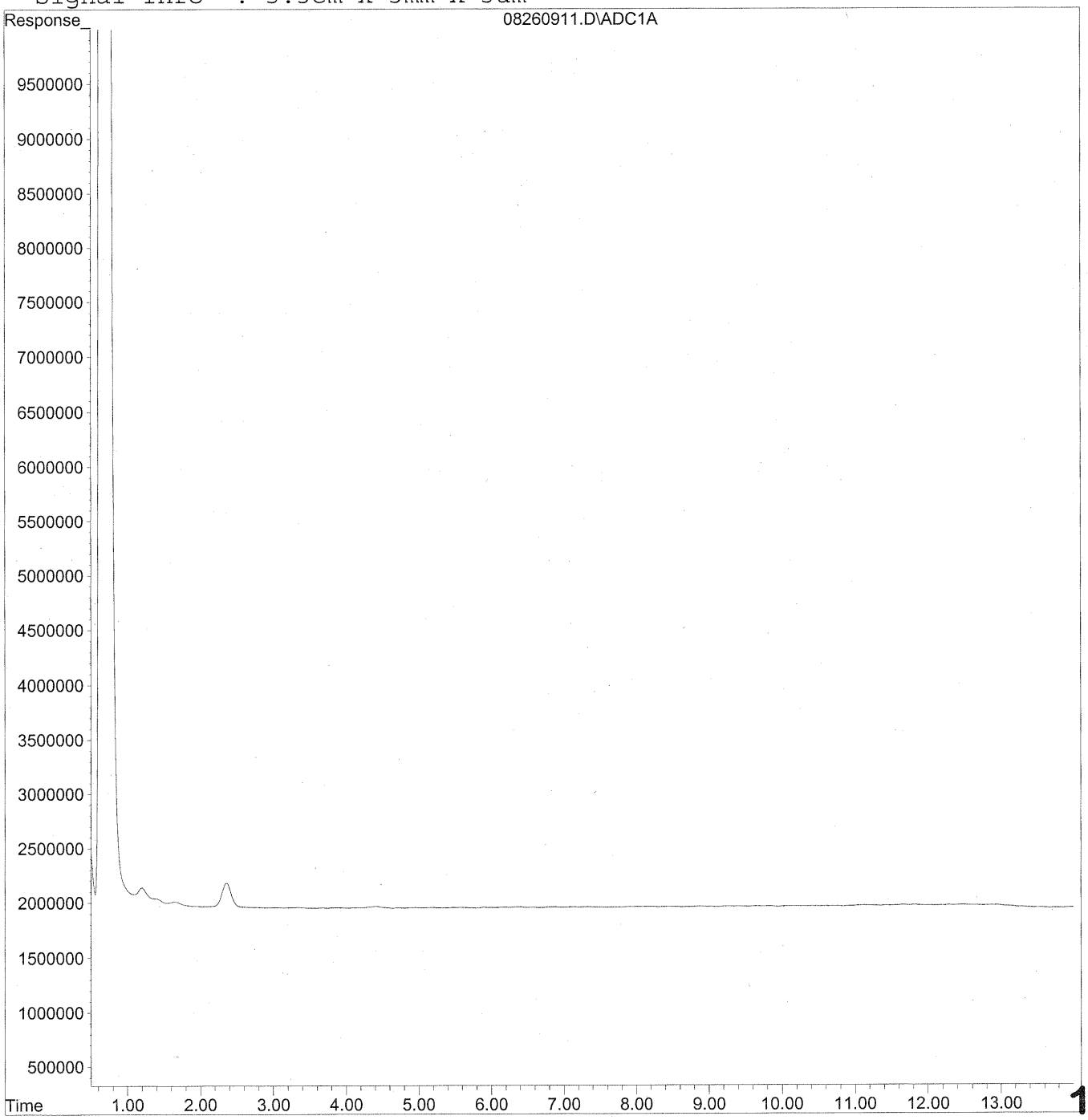
08270909.D TO110709.M Sun Aug 30 13:40:09 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260911.D Vial: 11
Acq On : 26 Aug 2009 7:35 pm Operator: HC
Sample : P0902942-006 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:51 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 : Thu Aug 27 07:35:56 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



149

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\26\08260911.D Vial: 11
Acq On : 26 Aug 2009 7:35 pm Operator: HC
Sample : P0902942-006 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:51 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 : Thu Aug 27 07:35:56 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

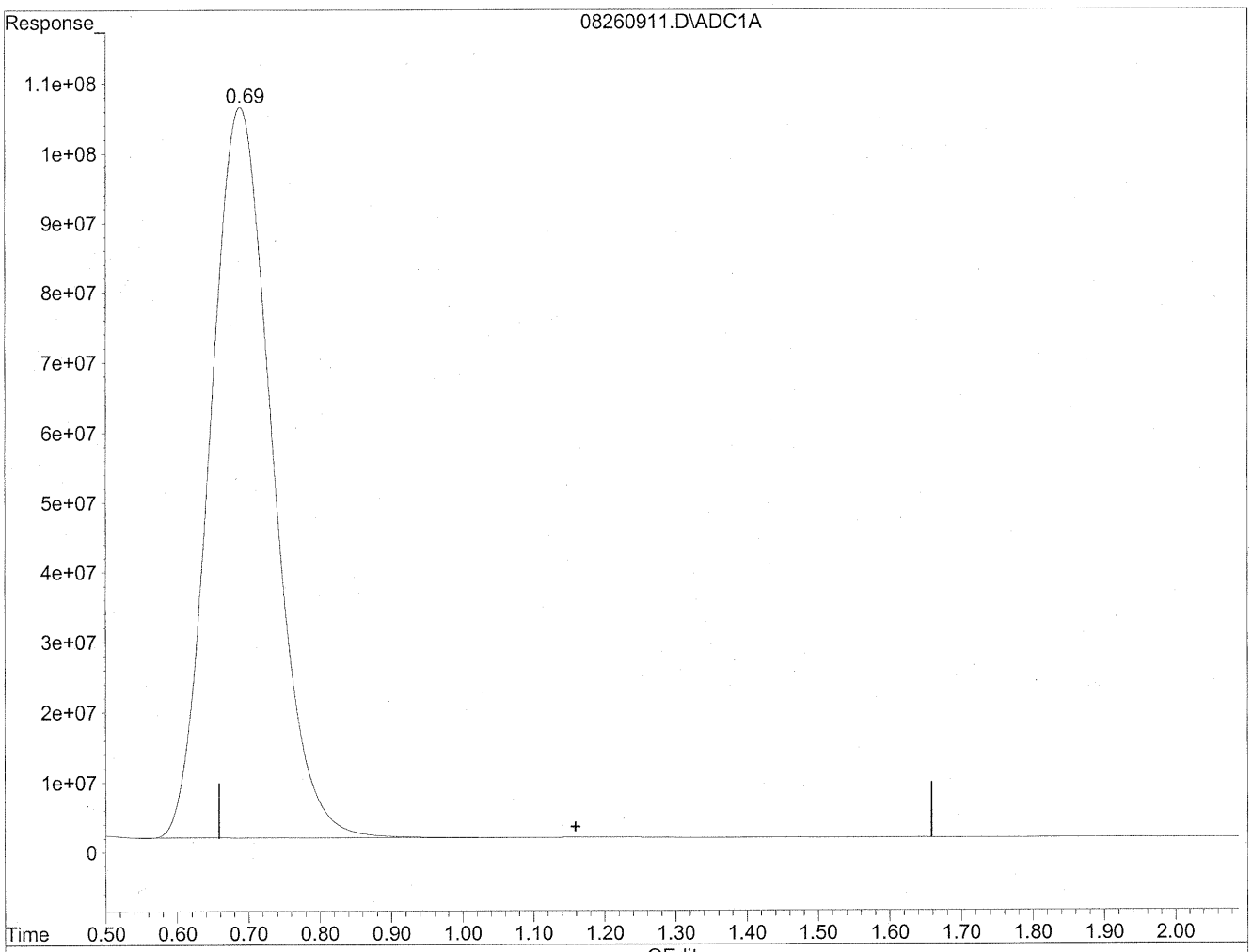
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	0.00	0	N.D.	ng/mld
2)	Acetaldehyde	0.00	0	N.D.	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/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260911.D Vial: 11
Acq On : 26 Aug 2009 7:35 pm Operator: HC
Sample : P0902942-006 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

0.69min 34365.758ng/ml

response 6308913851

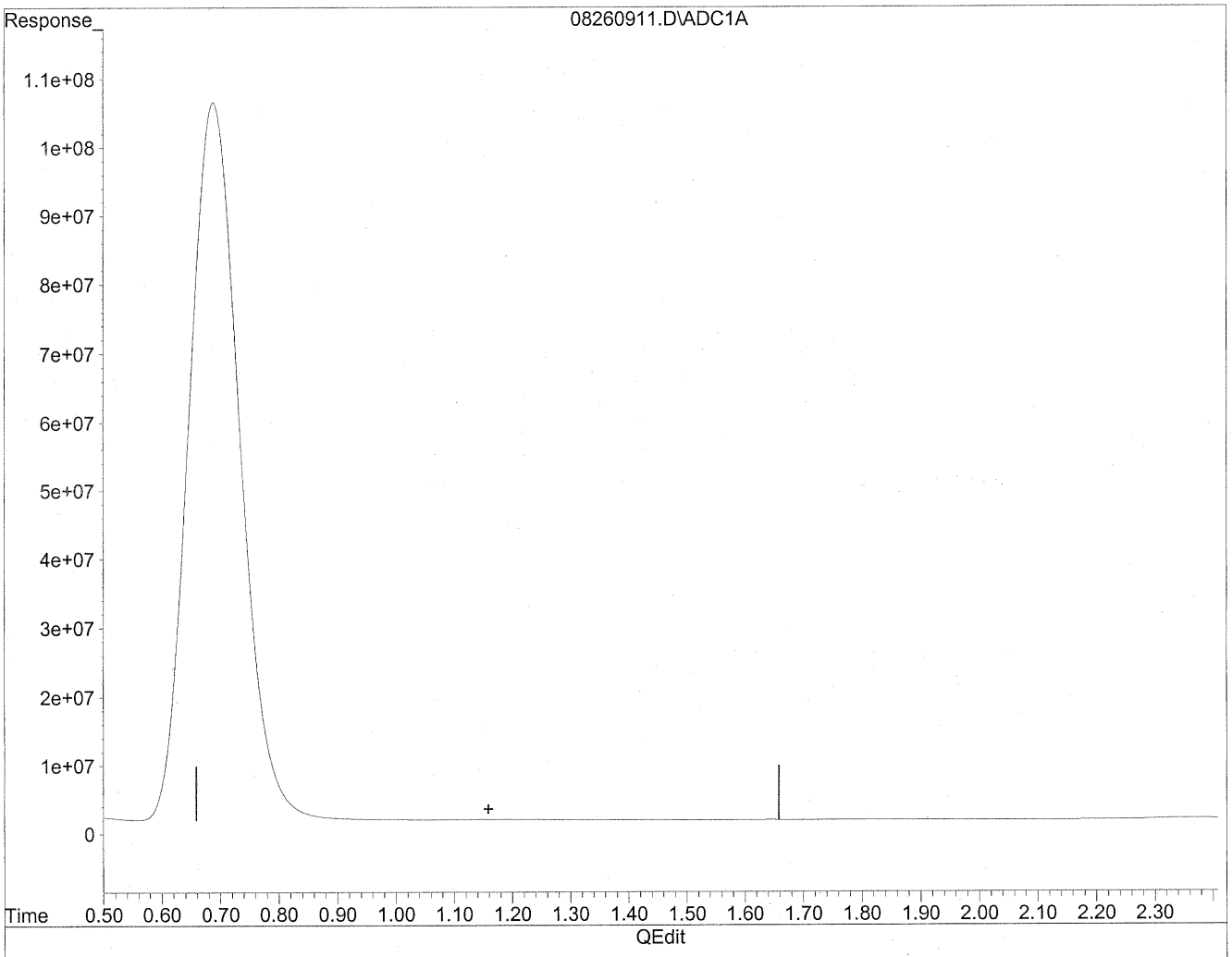
(+) = Expected Retention Time

08260911.D TO110709.M Sat Aug 29 16:50:58 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260911.D Vial: 11
Acq On : 26 Aug 2009 7:35 pm Operator: HC
Sample : P0902942-006 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:37 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

0.00min 0.000ng/ml d

response 0

HL 8/30/09 WP

8/31/09

152

(+) = Expected Retention Time

08260911.D TO110709.M Sat Aug 29 16:51:00 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: P0902942
CAS Sample ID: P090826-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: 08/26/09
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.

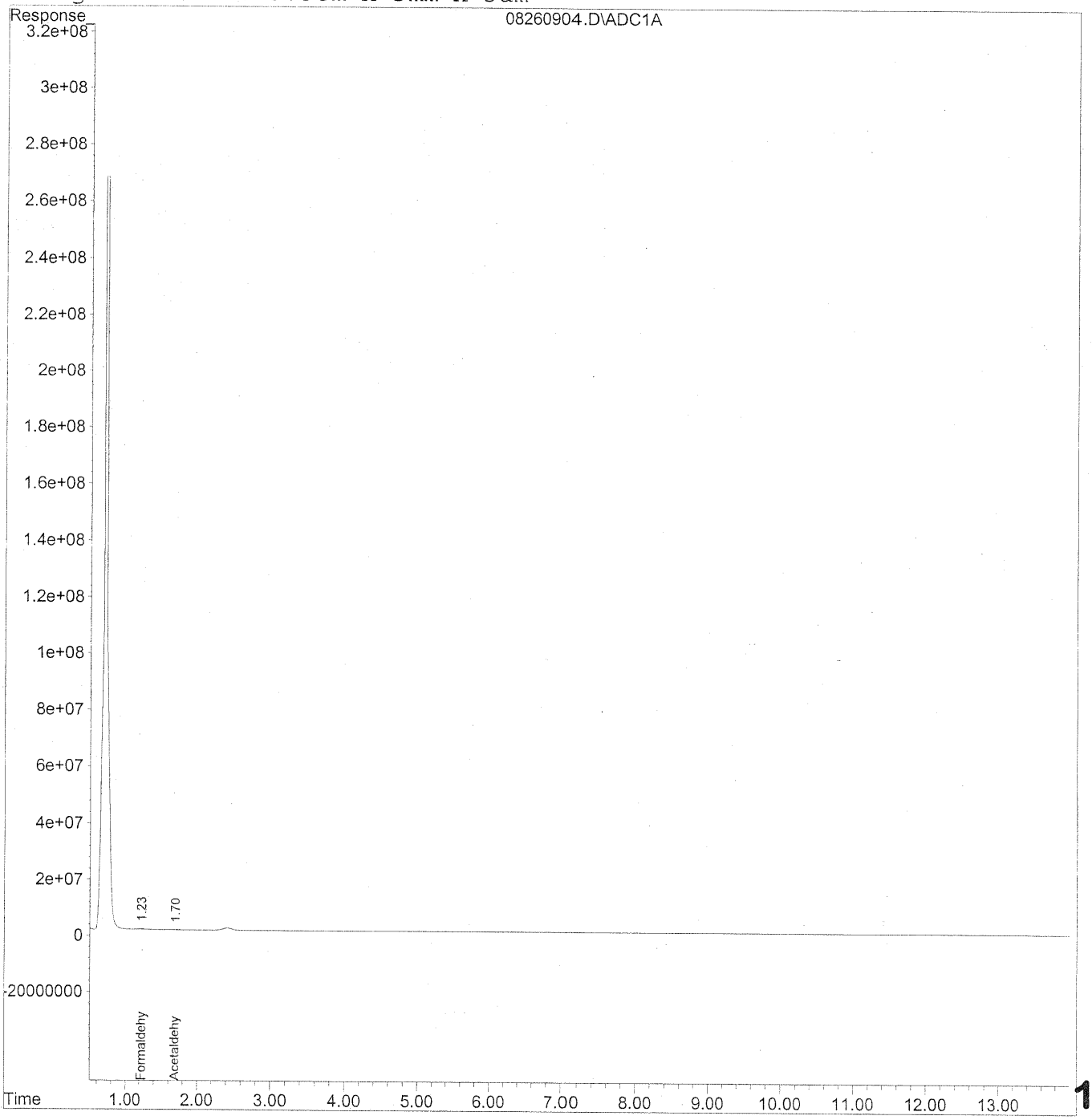
Verified By: RC Date: 9/11/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260904.D Vial: 4
Acq On : 26 Aug 2009 5:50 pm Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:31 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 : Sat Aug 29 12:41:27 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\26\08260904.D Vial: 4
Acq On : 26 Aug 2009 5:50 pm Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:31 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 : Sat Aug 29 12:41:27 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

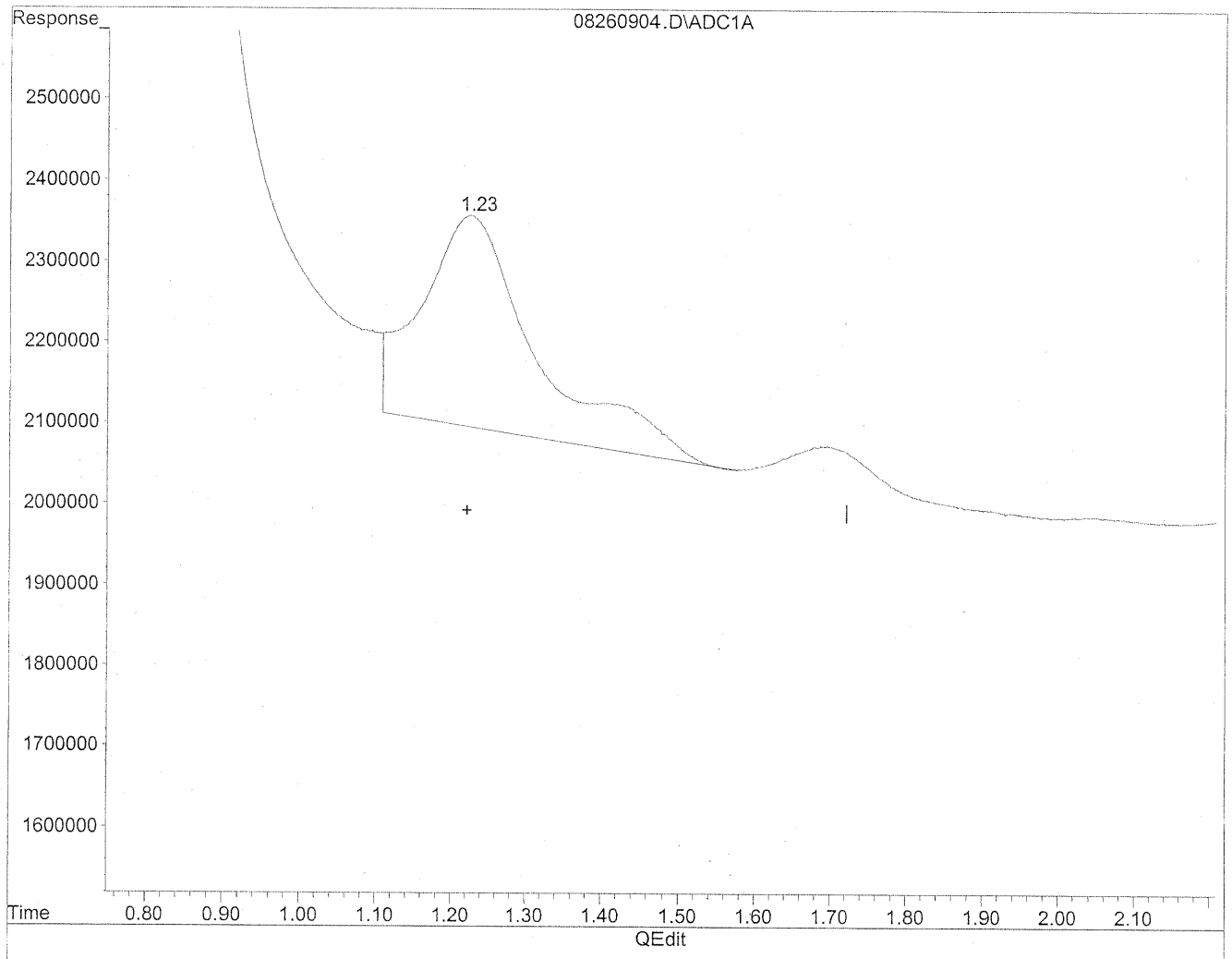
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.23	11782862	64.183	ng/mlm
2)	Acetaldehyde	1.70	3480548	24.821	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	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\26\08260904.D Vial: 4
Acq On : 26 Aug 2009 5:50 pm Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:30 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 : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.23min 151.780ng/ml

response 27863991

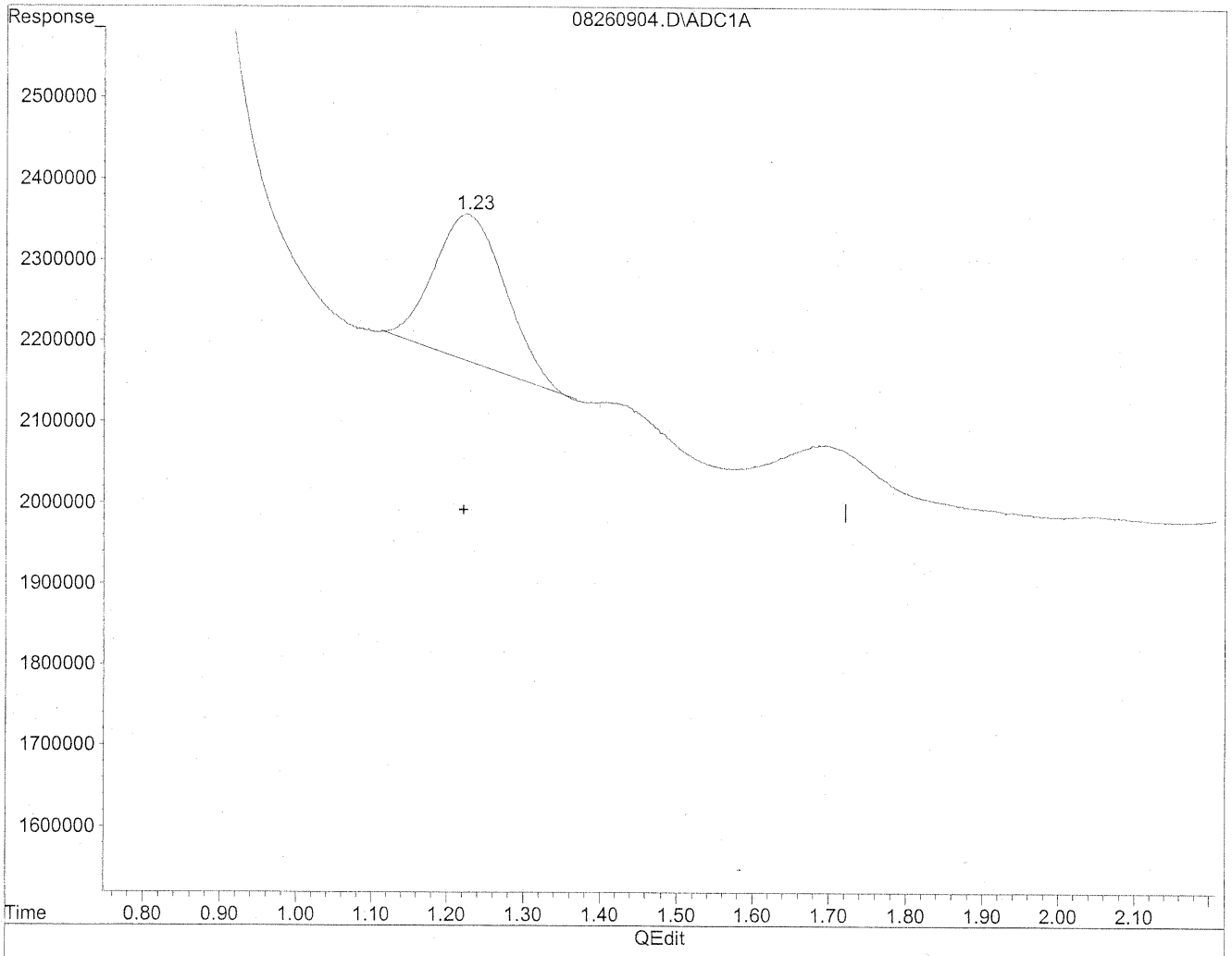
(+) = Expected Retention Time

08260904.D TO110709.M Sat Aug 29 16:31:08 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260904.D Vial: 4
Acq On : 26 Aug 2009 5:50 pm Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:30 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 : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.23min 64.183ng/ml m

response 11782862

tic 8/29/09 LC

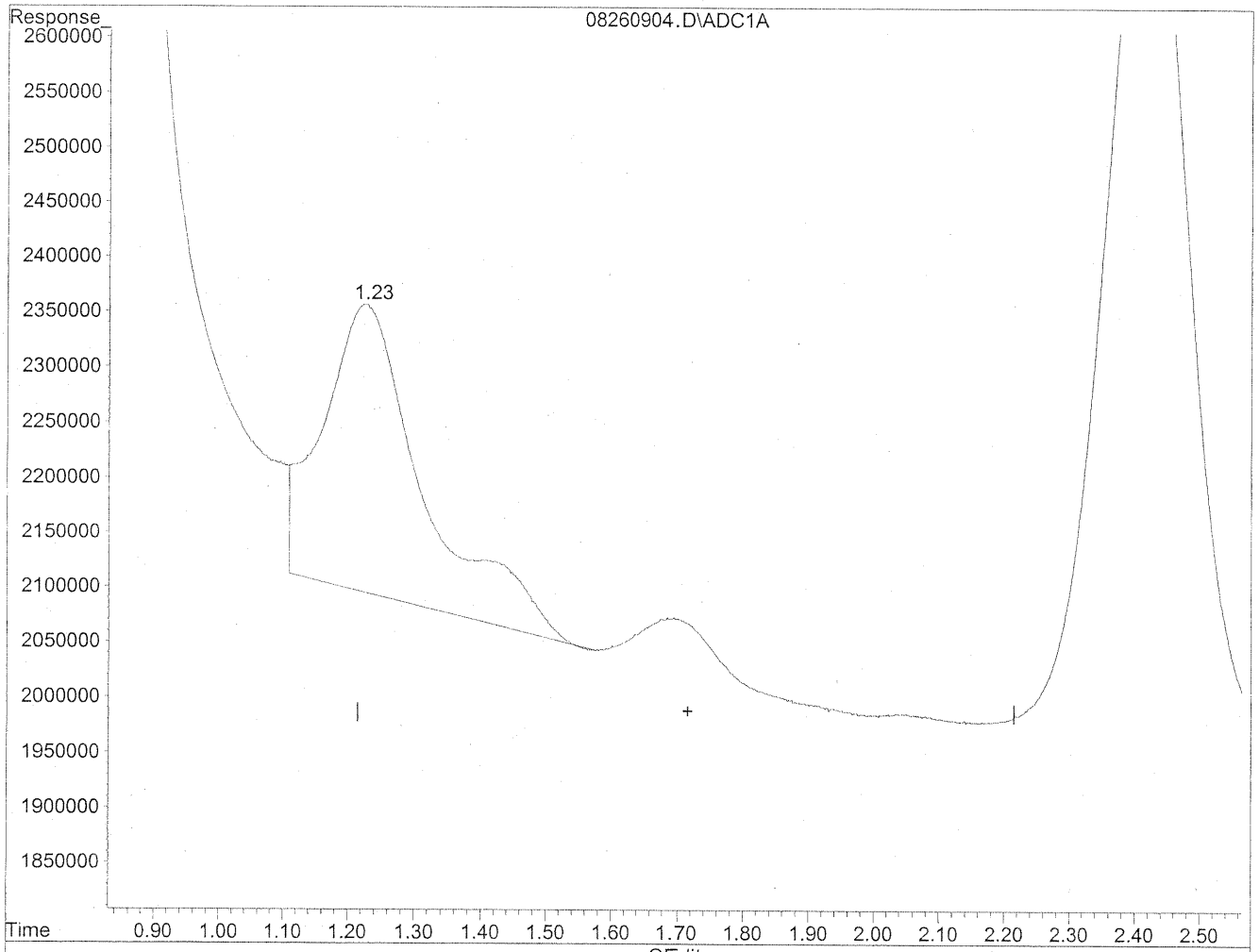
(+) = Expected Retention Time

08260904.D TO110709.M Sat Aug 29 16:31:15 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260904.D Vial: 4
Acq On : 26 Aug 2009 5:50 pm Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:30 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 : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.23min 198.711ng/ml

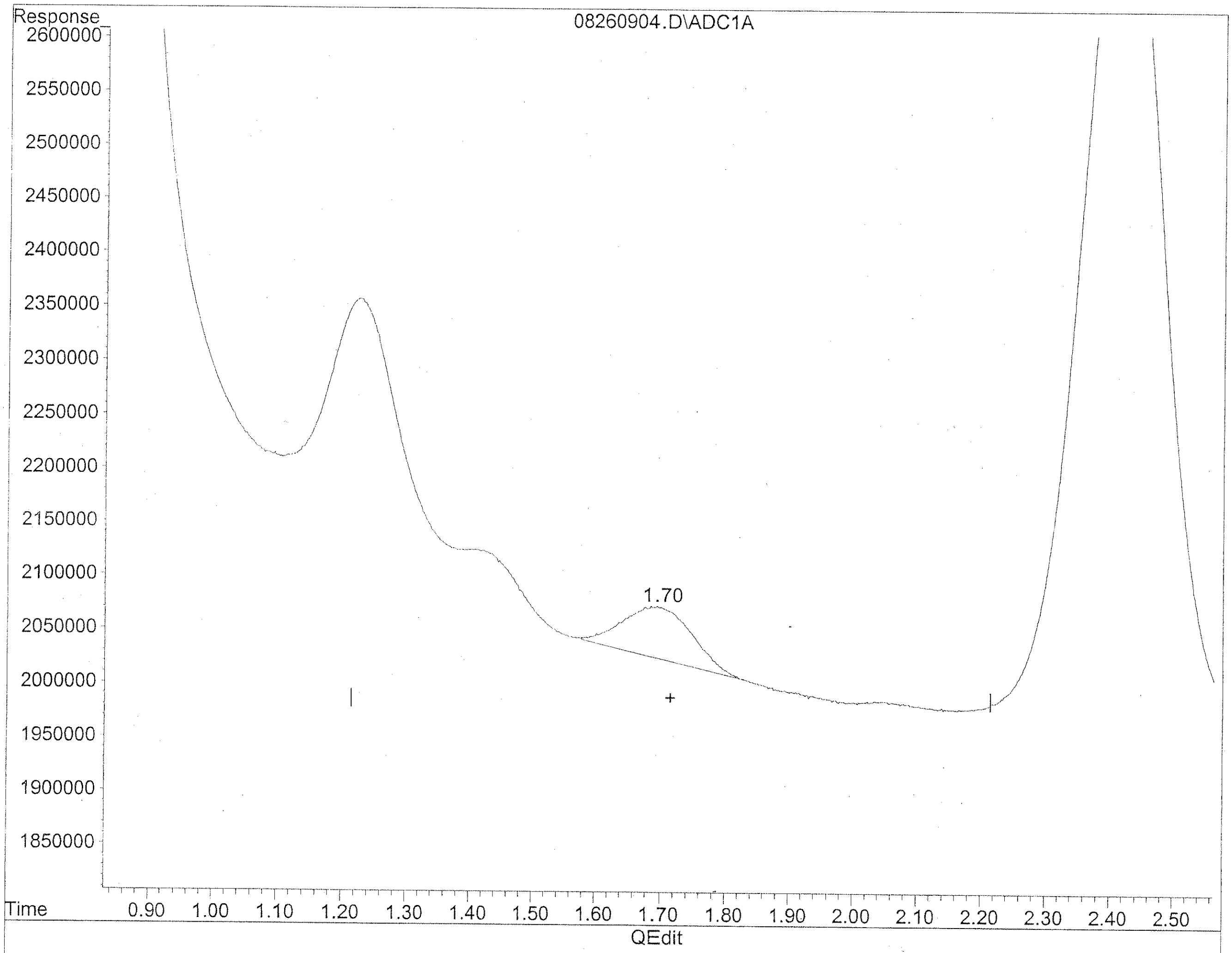
response 27863991

(+) = Expected Retention Time

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2009_08\26\08260904.D        Vial: 4
Acq On    : 26 Aug 2009   5:50 pm                      Operator: HC
Sample    : MB front lot 5855/5994 1.0ml               Inst    : LC 01
Misc      :                                            Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Aug 29 16:30 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     : Sat Aug 29 12:41:27 2009
Response via    : Multiple Level Calibration
```



(2) Acetaldehyde

1.70min 24.821ng/ml m

response 3480548

(+) = Expected Retention Time

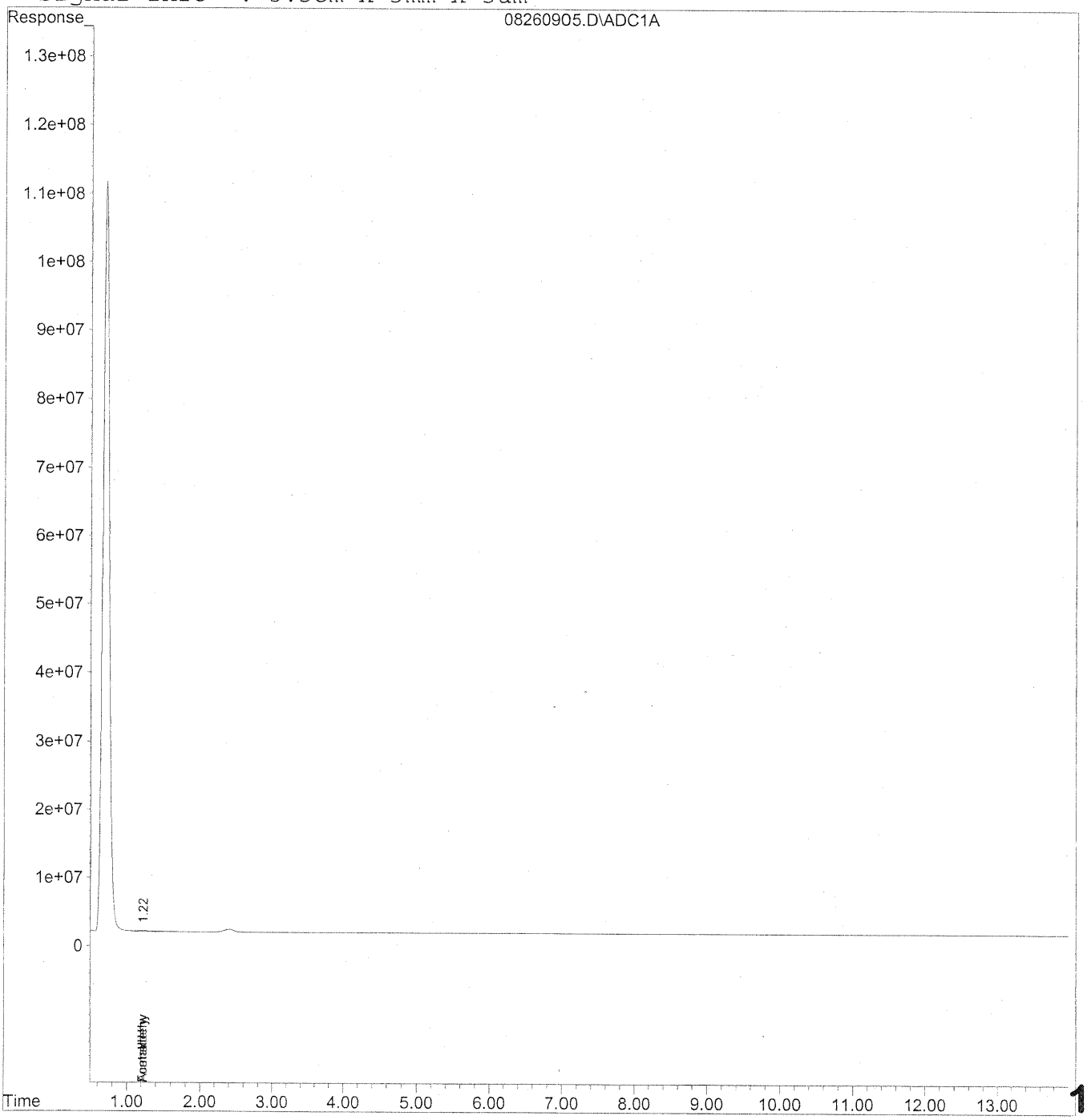
08260904.D TO110709.M Sat Aug 29 16:31:30 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260905.D Vial: 5
Acq On : 26 Aug 2009 6:05 pm Operator: HC
Sample : MB back lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:32 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 : Sat Aug 29 12:41:27 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\26\08260905.D Vial: 5
Acq On : 26 Aug 2009 6:05 pm Operator: HC
Sample : MB back lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:32 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 : Sat Aug 29 12:41:27 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	10945941	59.624	ng/ml
2)	Acetaldehyde	1.22f	10945941	78.061	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/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

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: P0902942
CAS Sample ID: P090827-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: 08/27/09
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	

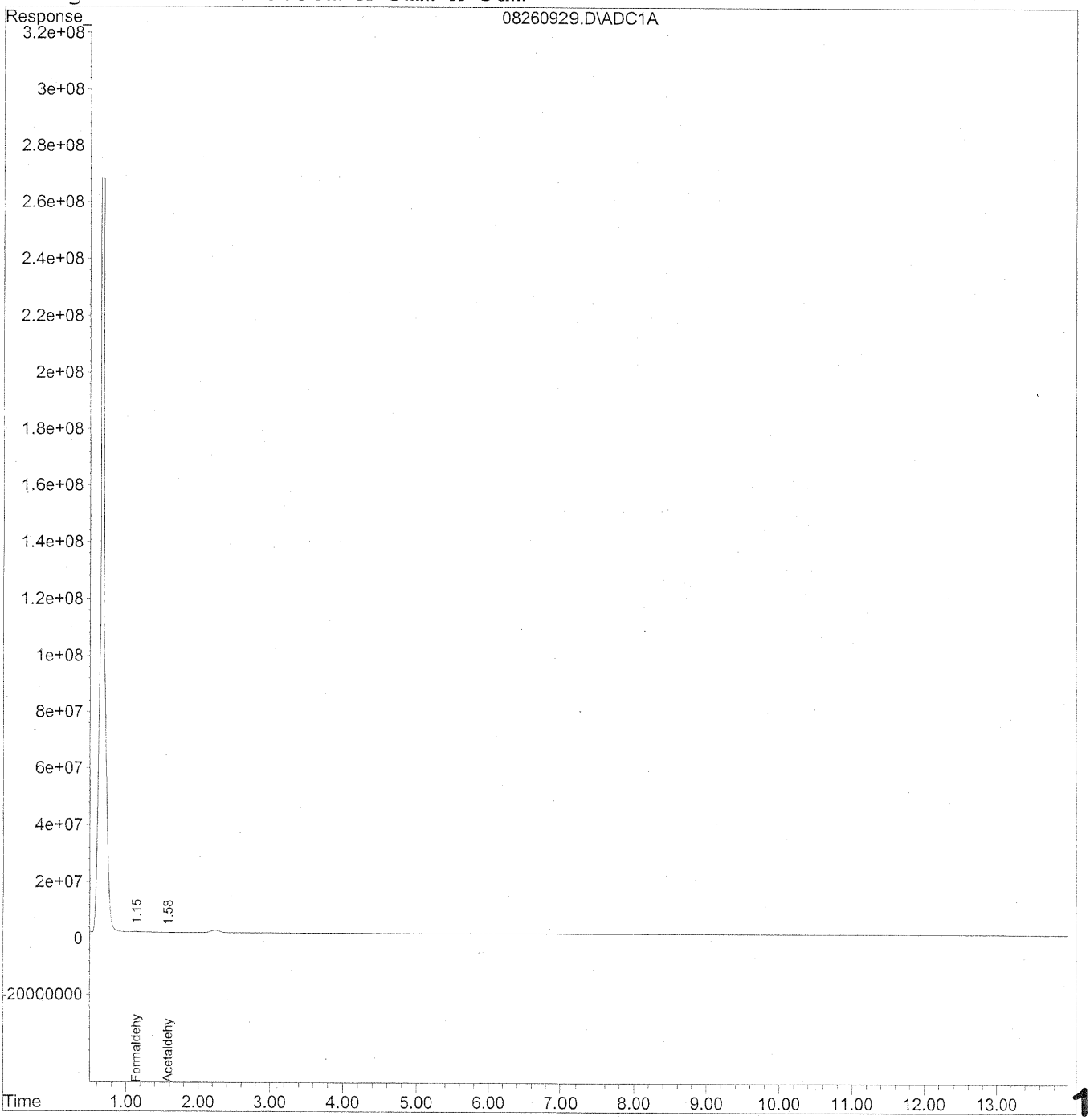
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
NA = Not applicable.

Verified By: Re Date: 9/11/09

Data File : J:\LC01\DATA\TO11\2009_08\26\08260929.D Vial: 28
Acq On : 27 Aug 2009 12:06 am Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:35 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 : Sat Aug 29 16:33:38 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\26\08260929.D Vial: 28
Acq On : 27 Aug 2009 12:06 am Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:35 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 : Sat Aug 29 16:33:38 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

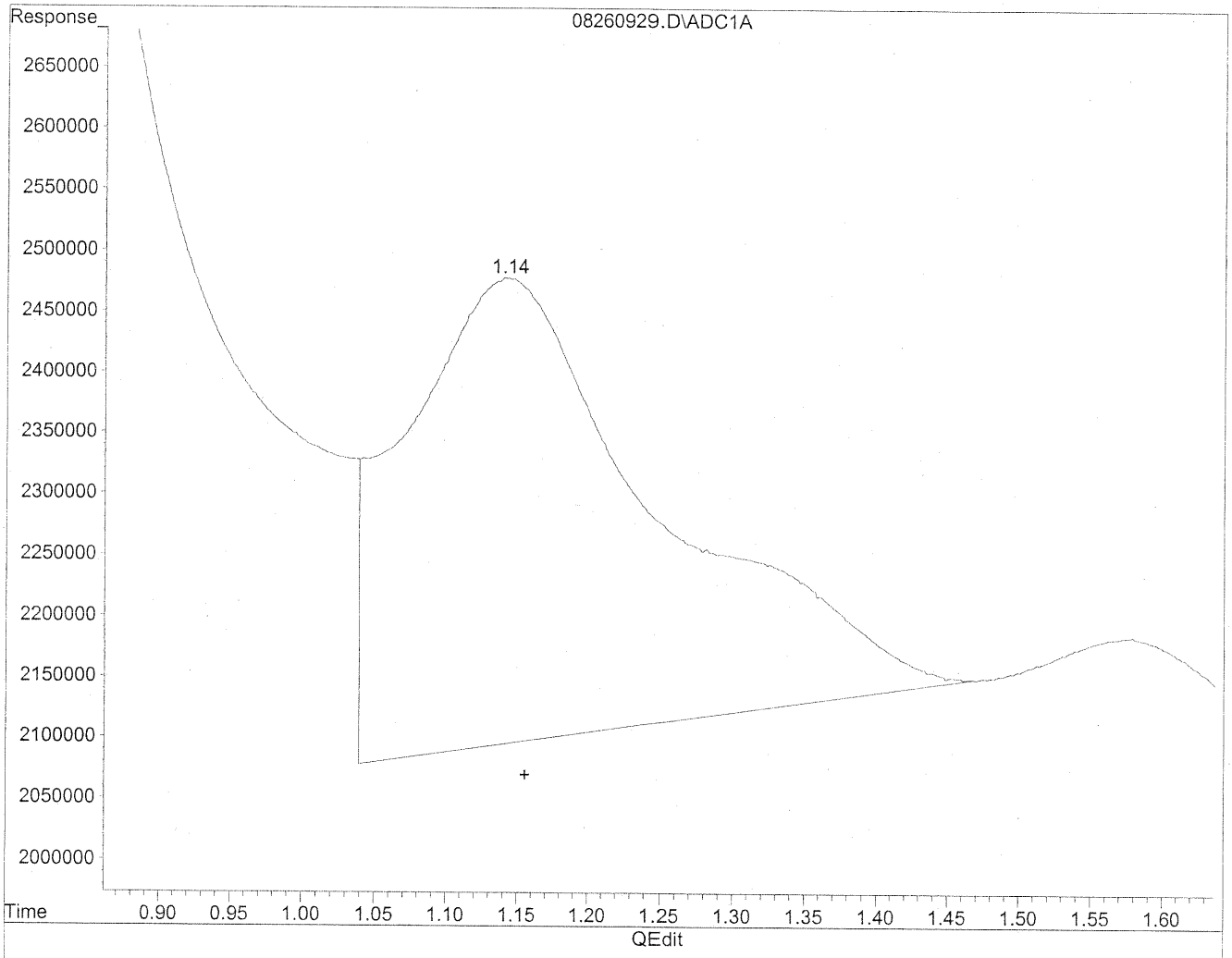
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.15	11602069	63.199	ng/mlm
2)	Acetaldehyde	1.58	4276541	30.498	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\26\08260929.D Vial: 28
Acq On : 27 Aug 2009 12:06 am Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:34 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.14min 257.970ng/ml

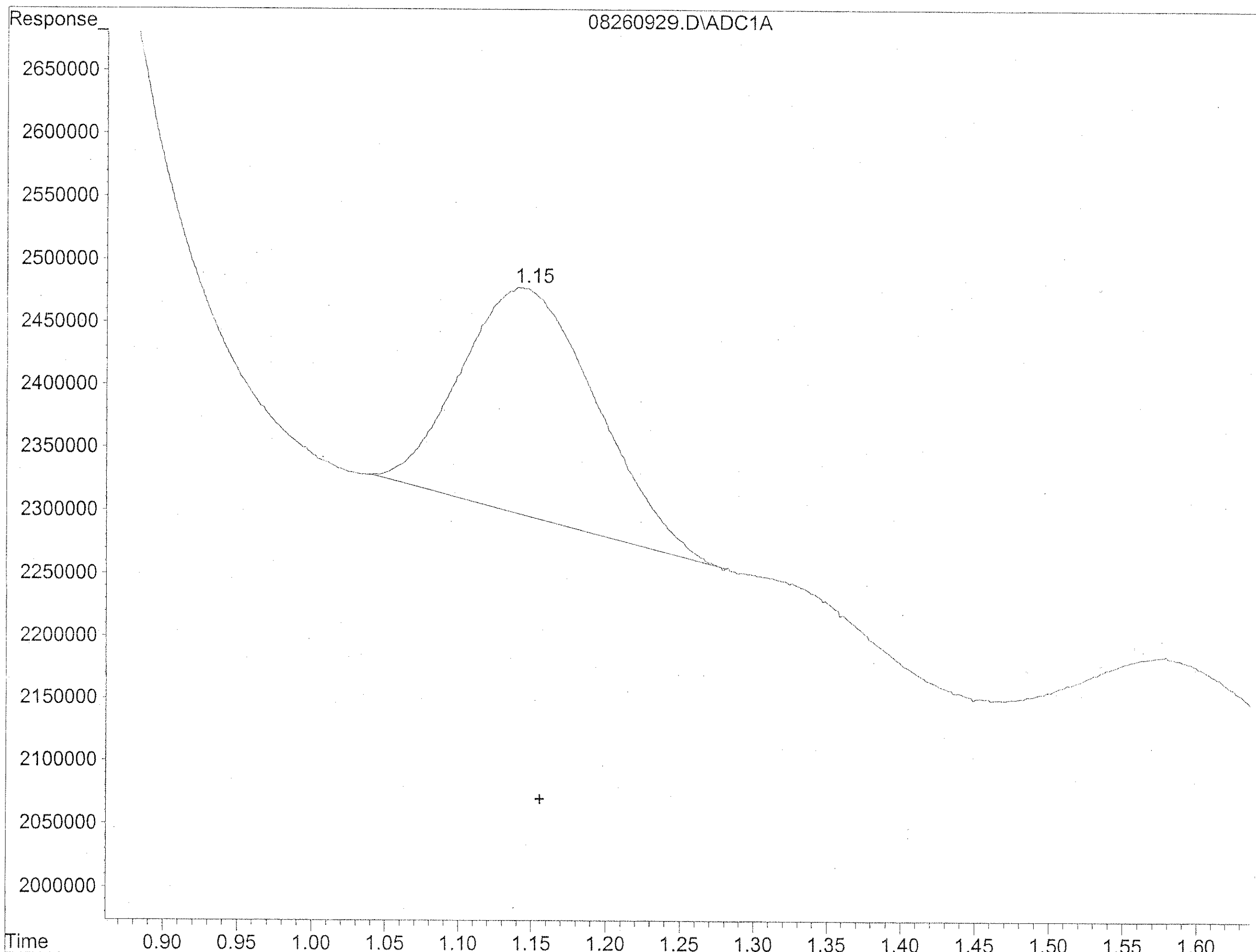
response 47358406

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260929.D Vial: 28
Acq On : 27 Aug 2009 12:06 am Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:34 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.15min 63.199ng/ml m

response 11602069

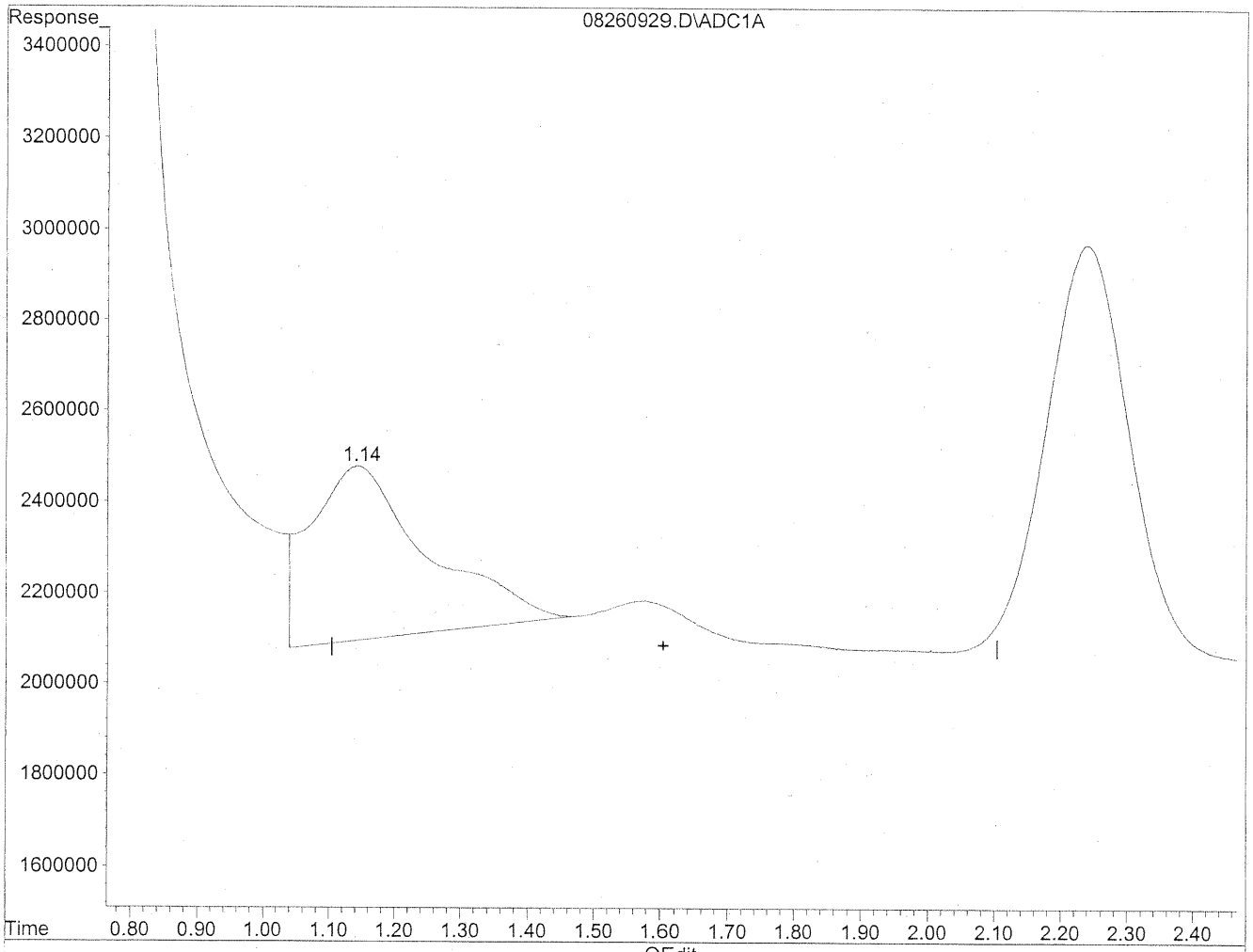
(+) = Expected Retention Time

08260929.D TO110709.M Sat Aug 29 16:35:11 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260929.D Vial: 28
Acq On : 27 Aug 2009 12:06 am Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:34 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.14min 337.735ng/ml

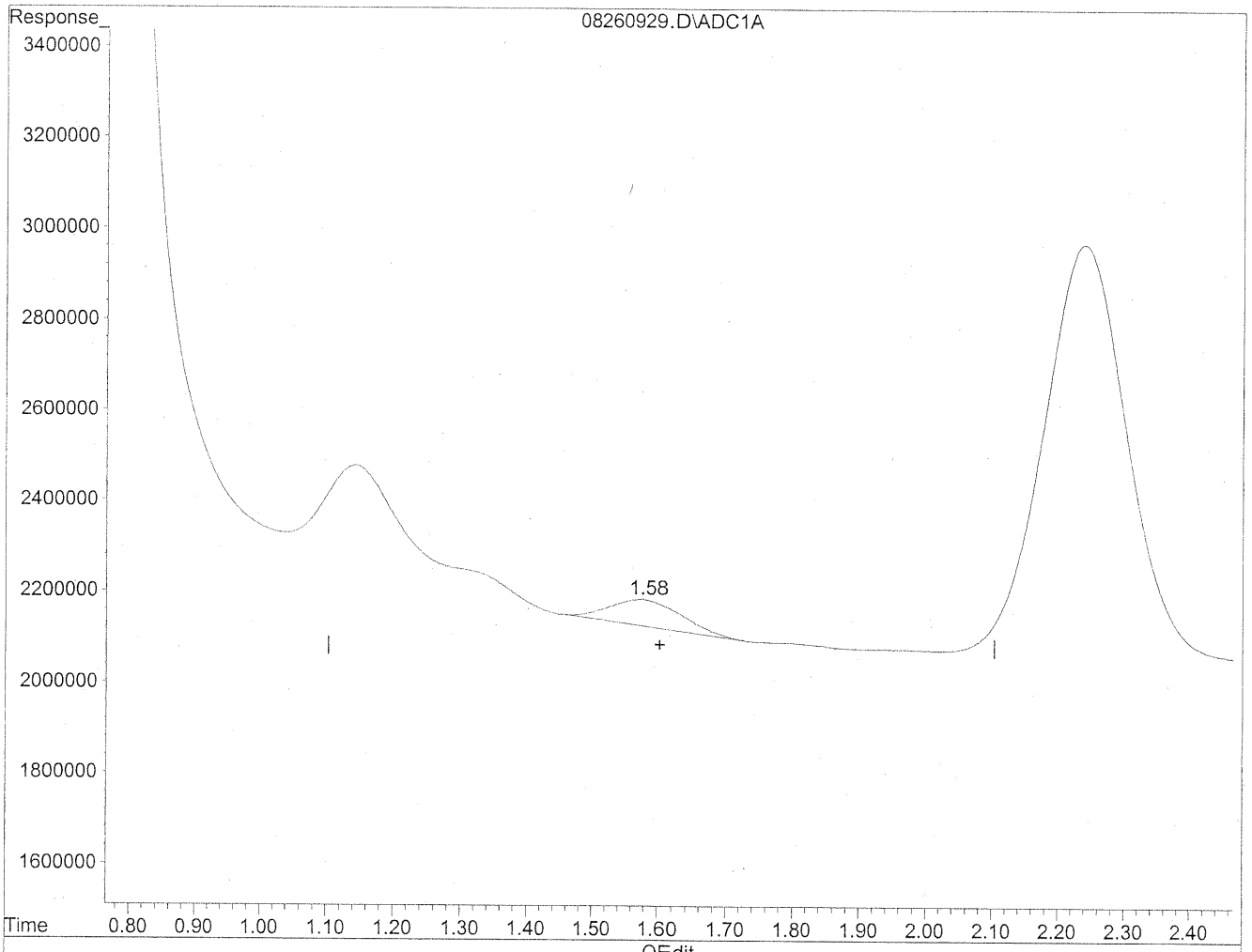
response 47358406

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260929.D Vial: 28
Acq On : 27 Aug 2009 12:06 am Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:34 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.58min 30.498ng/ml m

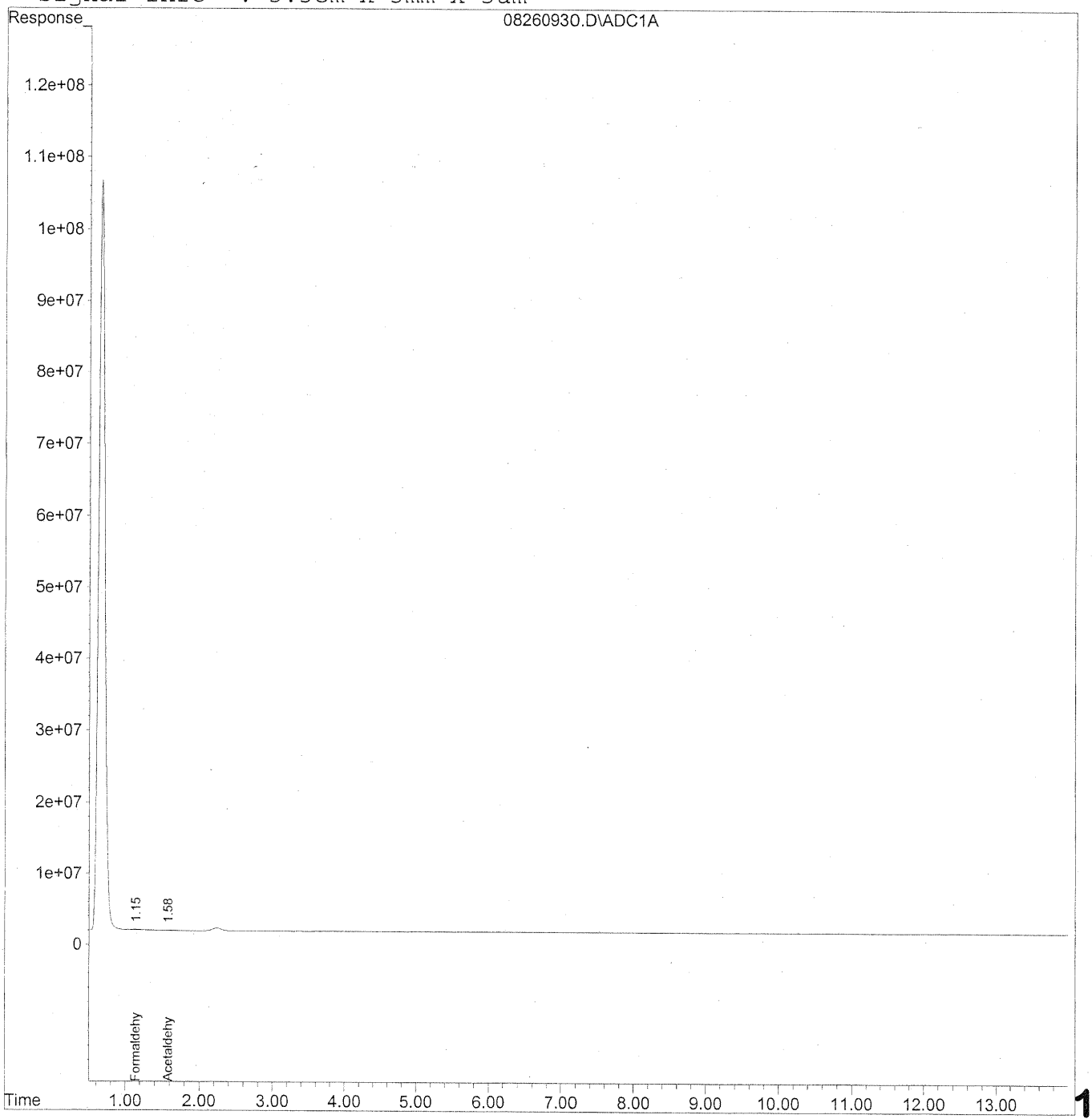
response 4276541

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_08\26\08260930.D Vial: 29
Acq On : 27 Aug 2009 12:21 am Operator: HC
Sample : MB back lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:36 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 : Sat Aug 29 16:33:38 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\26\08260930.D Vial: 29
Acq On : 27 Aug 2009 12:21 am Operator: HC
Sample : MB back lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:36 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 : Sat Aug 29 16:33:38 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

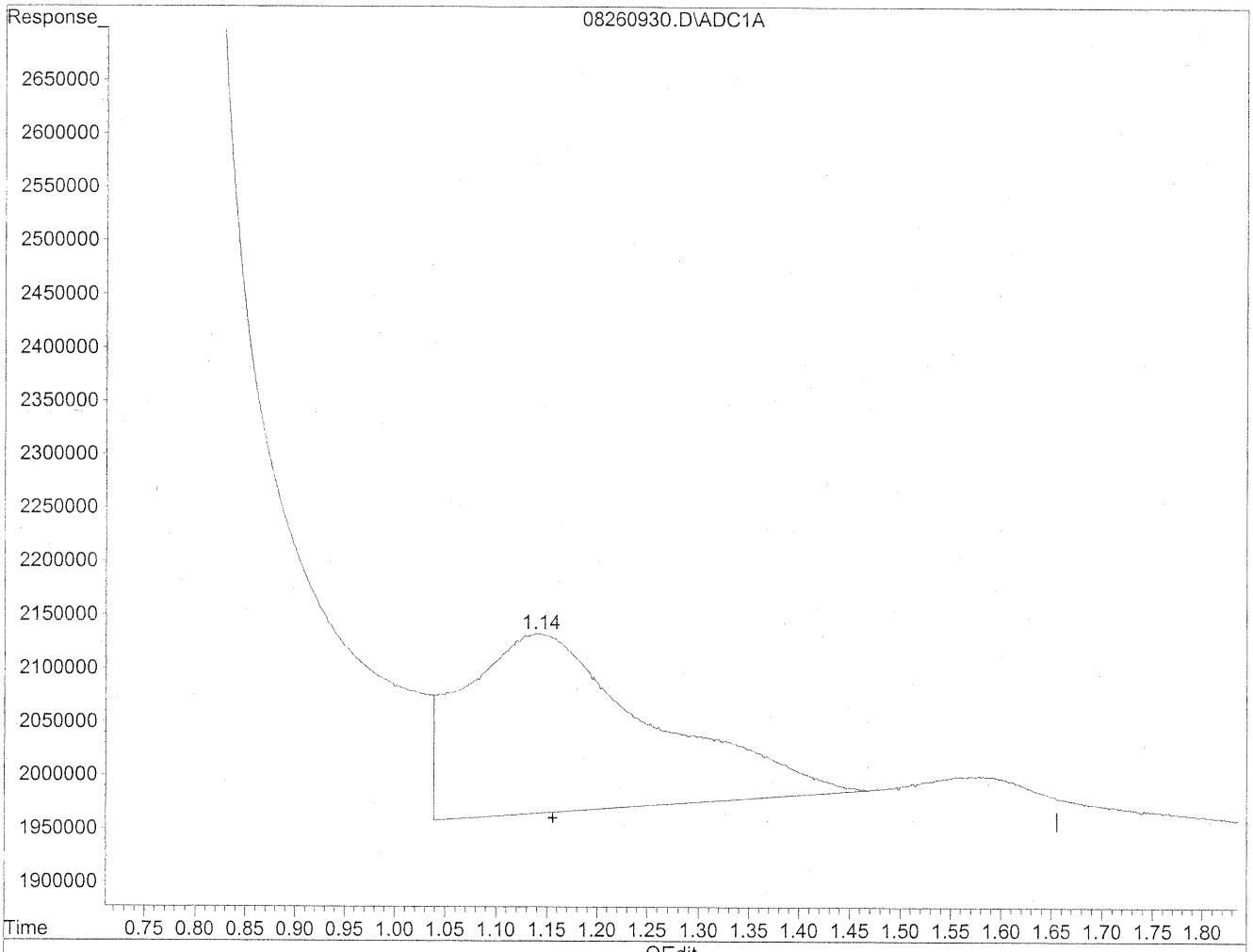
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.15	4503102	24.529	ng/mlm
2)	Acetaldehyde	1.58	1370787	9.776	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\26\08260930.D Vial: 29
Acq On : 27 Aug 2009 12:21 am Operator: HC
Sample : MB back lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:35 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.14min 118.486ng/ml

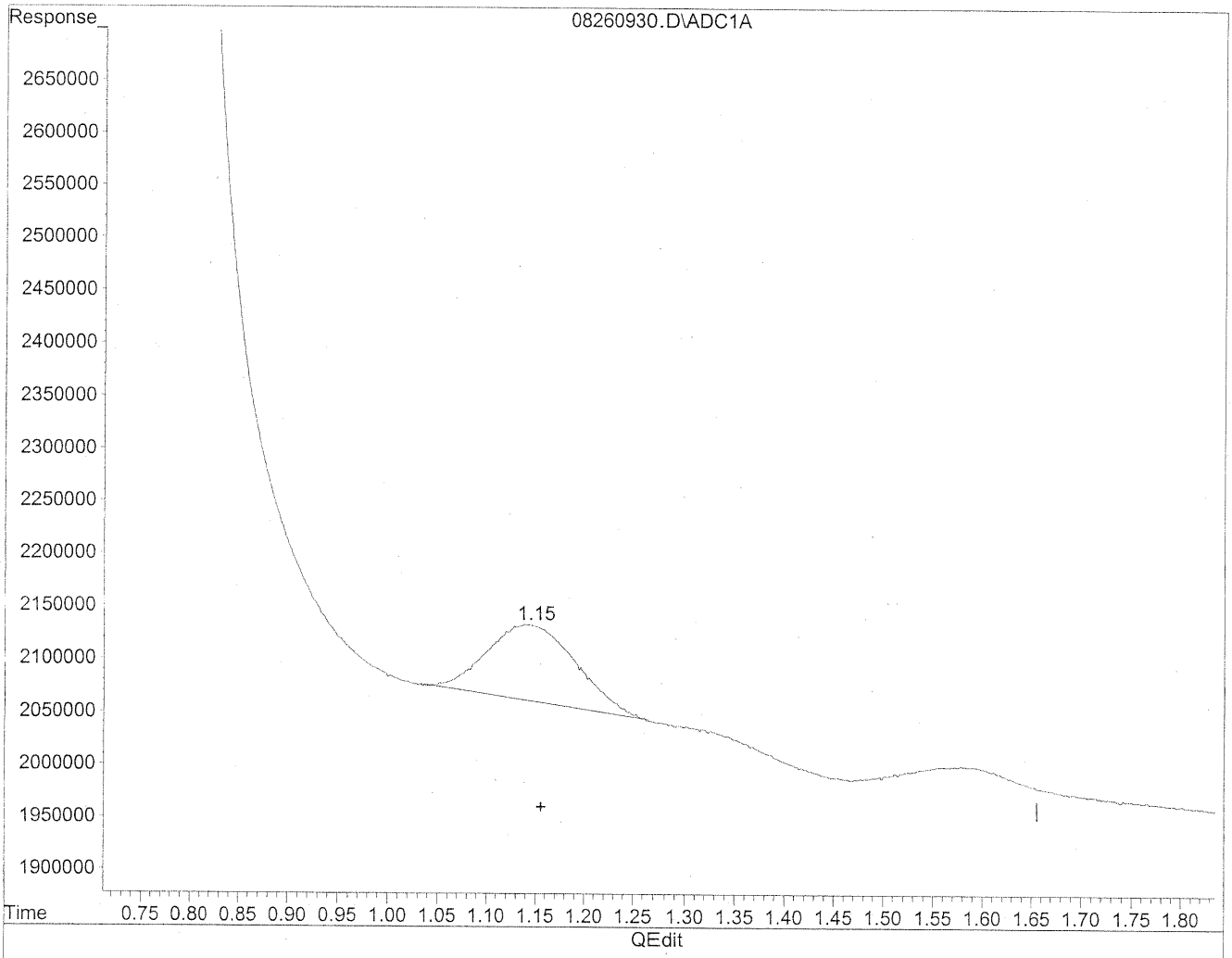
response 21751740

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260930.D Vial: 29
Acq On : 27 Aug 2009 12:21 am Operator: HC
Sample : MB back lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:35 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.15min 24.529ng/ml m

response 4503102

HC 8/29/09 LC

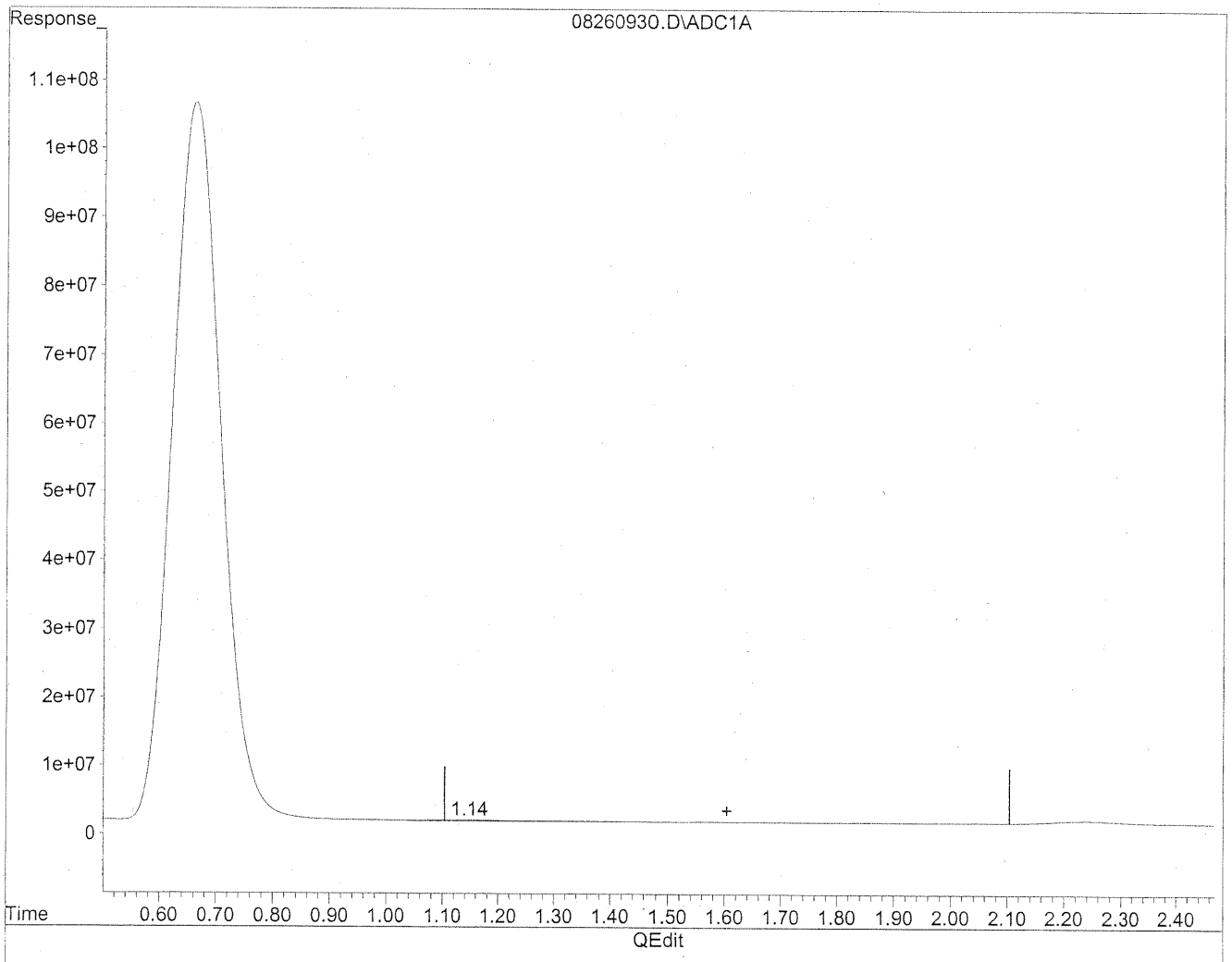
(+) = Expected Retention Time

08260930.D TO110709.M Sat Aug 29 16:35:54 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260930.D Vial: 29
Acq On : 27 Aug 2009 12:21 am Operator: HC
Sample : MB back lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:35 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.14min 155.122ng/ml

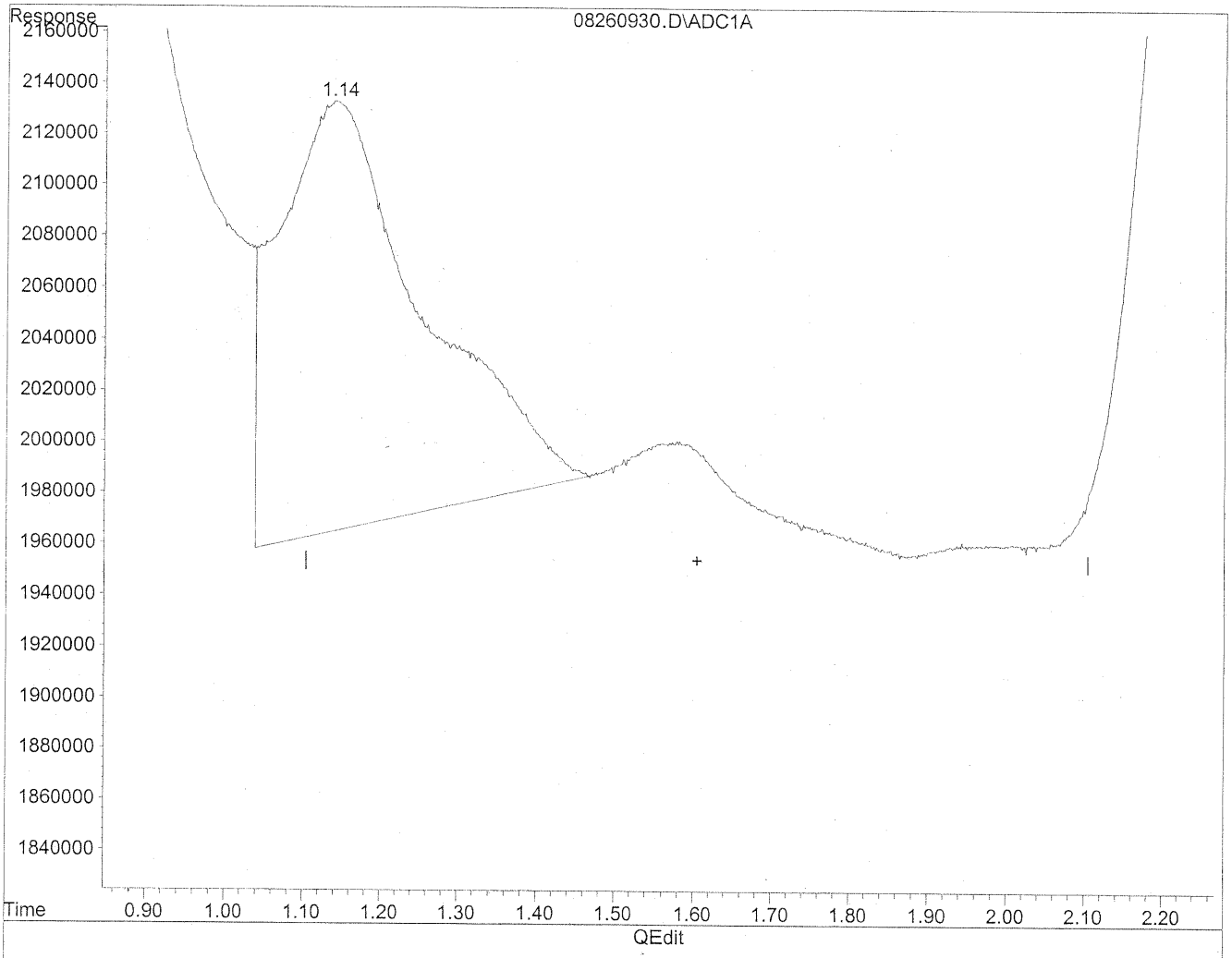
response 21751740

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260930.D Vial: 29
Acq On : 27 Aug 2009 12:21 am Operator: HC
Sample : MB back lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:35 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.14min 155.122ng/ml

response 21751740

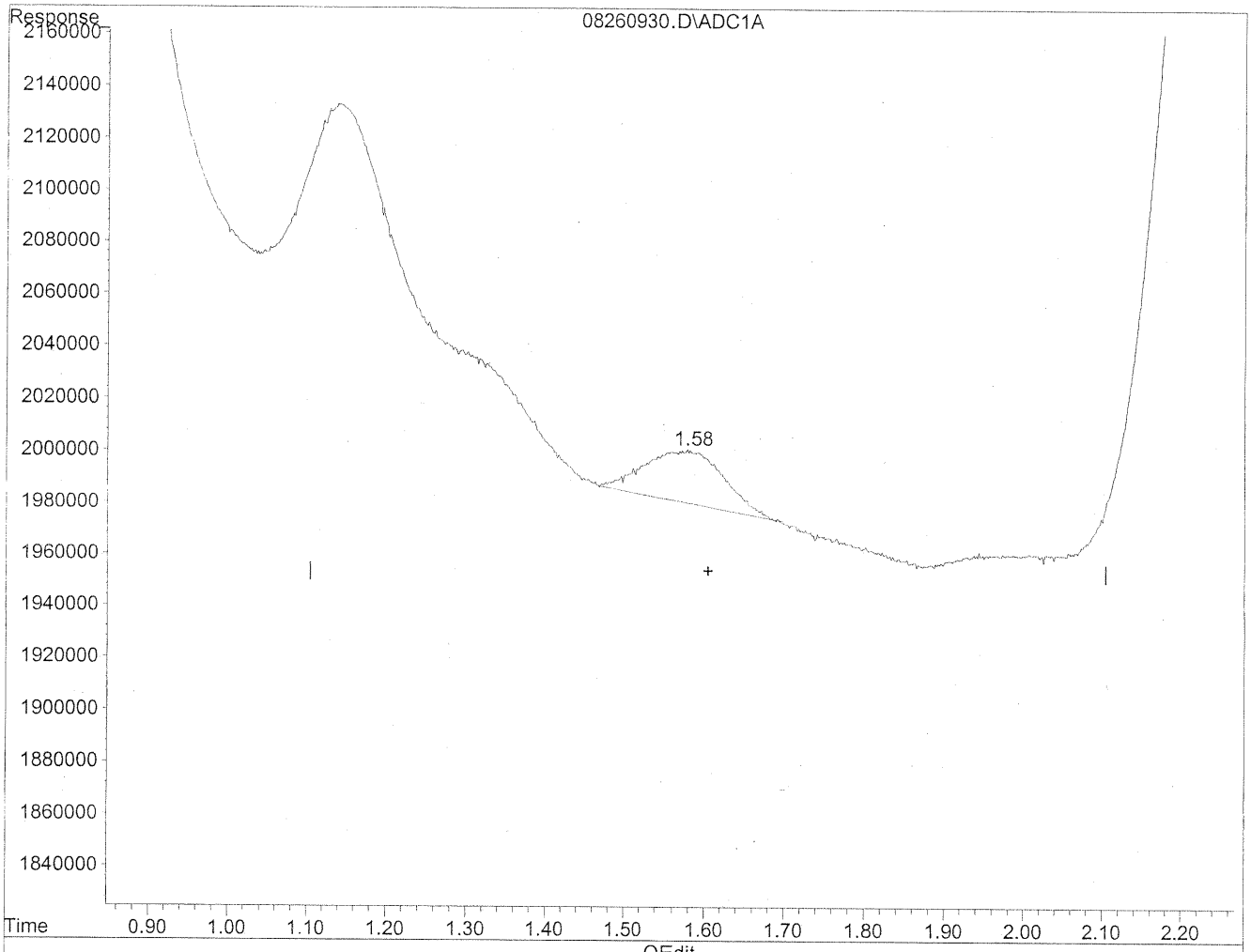
(+) = Expected Retention Time

08260930.D TO110709.M Sat Aug 29 16:36:03 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260930.D Vial: 29
Acq On : 27 Aug 2009 12:21 am Operator: HC
Sample : MB back lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:35 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.58min 9.776ng/ml m

response 1370787

(+) = Expected Retention Time

08260930.D TO110709.M Sat Aug 29 16:36:09 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **Method Blank (09:35)**
Client Project ID: **16512**

CAS Project ID: P0902942
CAS Sample ID: P090827-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: 08/27/09
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	< 200	NA	NA	NA	NA	**V**
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	< 200	NA	NA	NA	NA	**V**

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.
V = The continuing calibration verification standard was outside (biased low) the specified limits for this compound.

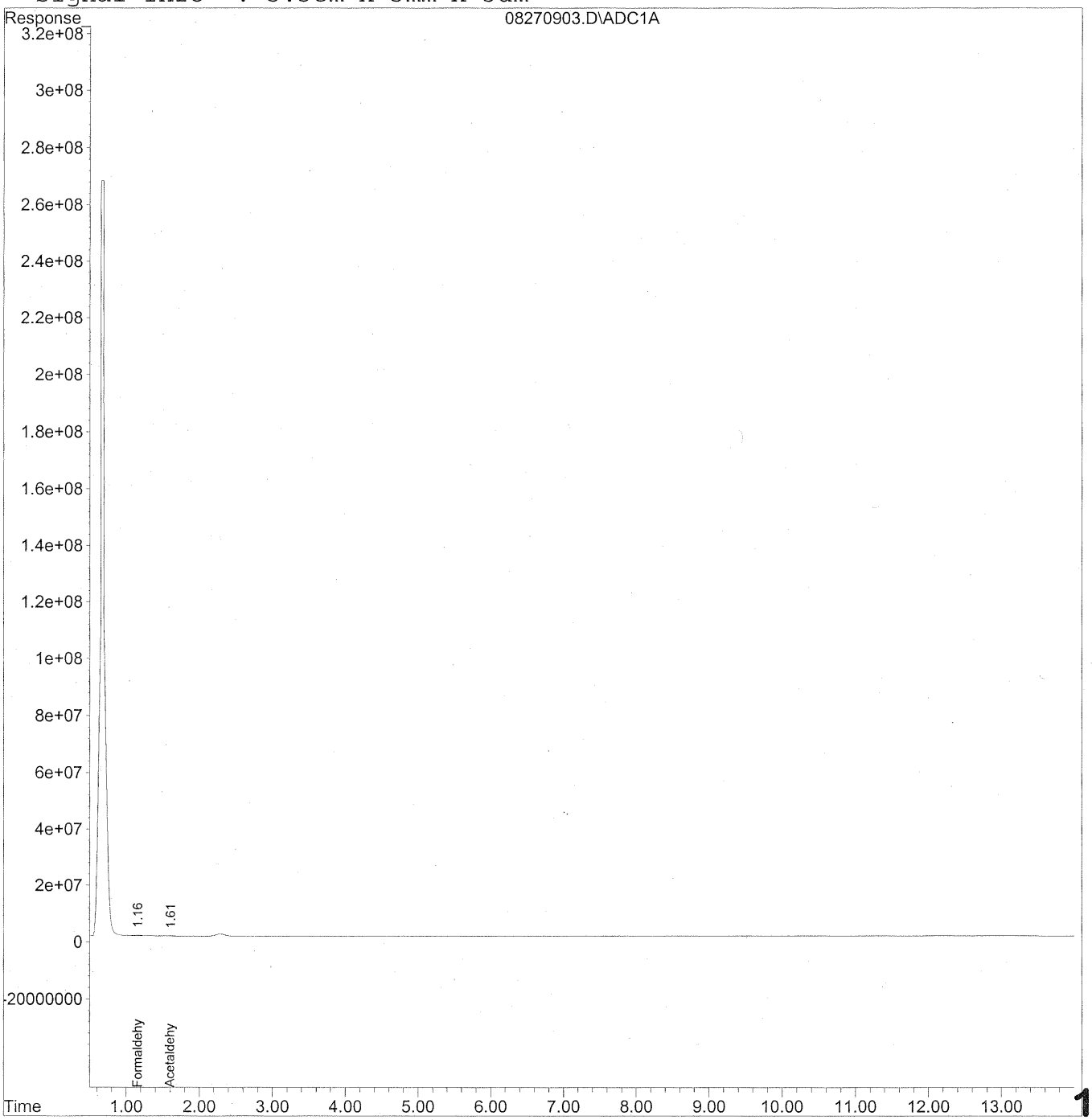
Verified By: Re. Date: 9/10/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270903.D Vial: 3
Acq On : 27 Aug 2009 9:35 am Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:43 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 : Sat Aug 29 16:33:38 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



177

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\27\08270903.D Vial: 3
Acq On : 27 Aug 2009 9:35 am Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:43 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 : Sat Aug 29 16:33:38 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

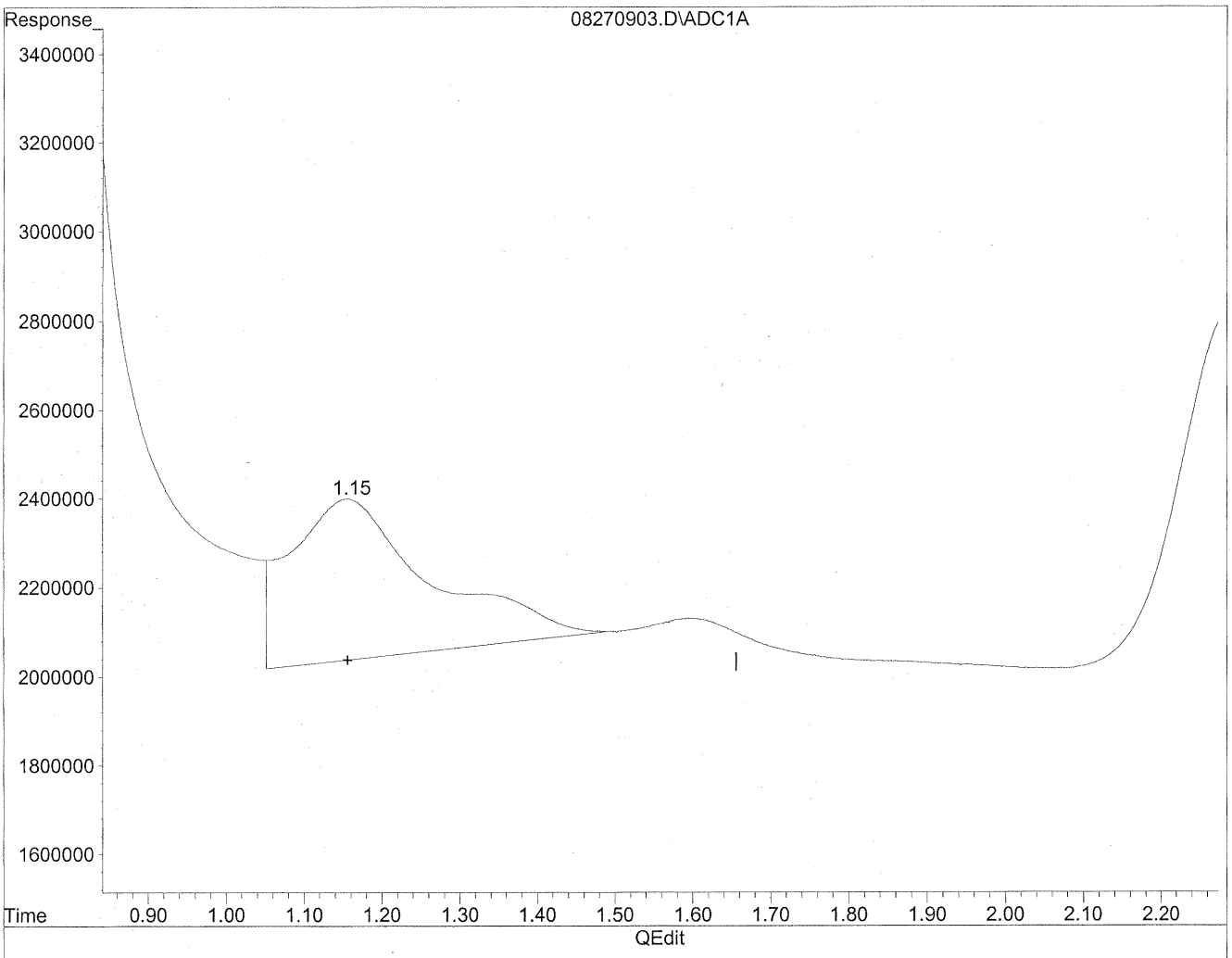
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
Target Compounds					
1)	Formaldehyde	1.16	11166974	60.828	ng/mlm
2)	Acetaldehyde	1.61	4012978	28.618	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\27\08270903.D Vial: 3
Acq On : 27 Aug 2009 9:35 am Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:42 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.16min 244.318ng/ml

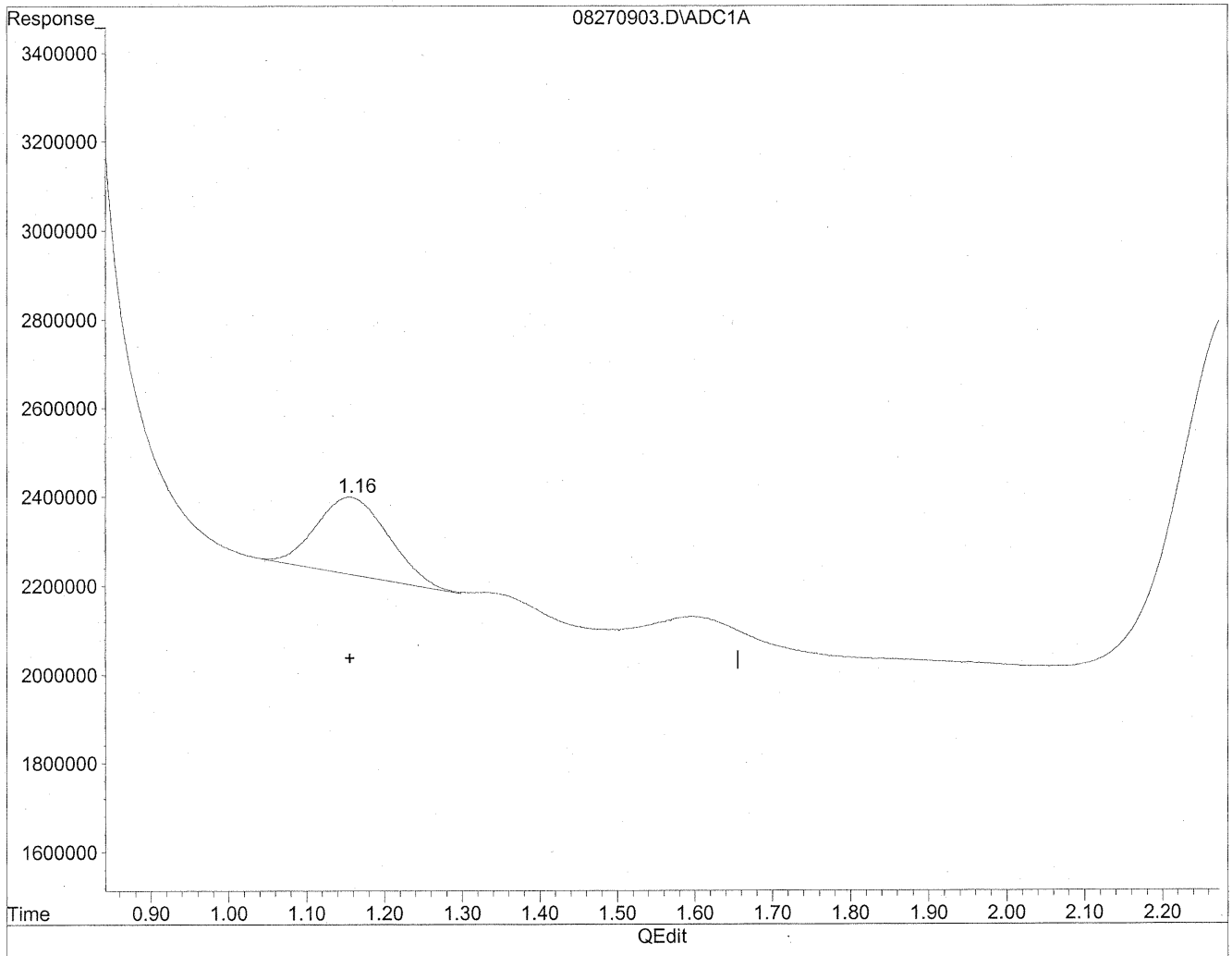
response 44852245

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270903.D Vial: 3
Acq On : 27 Aug 2009 9:35 am Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:42 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.16min 60.828ng/ml m

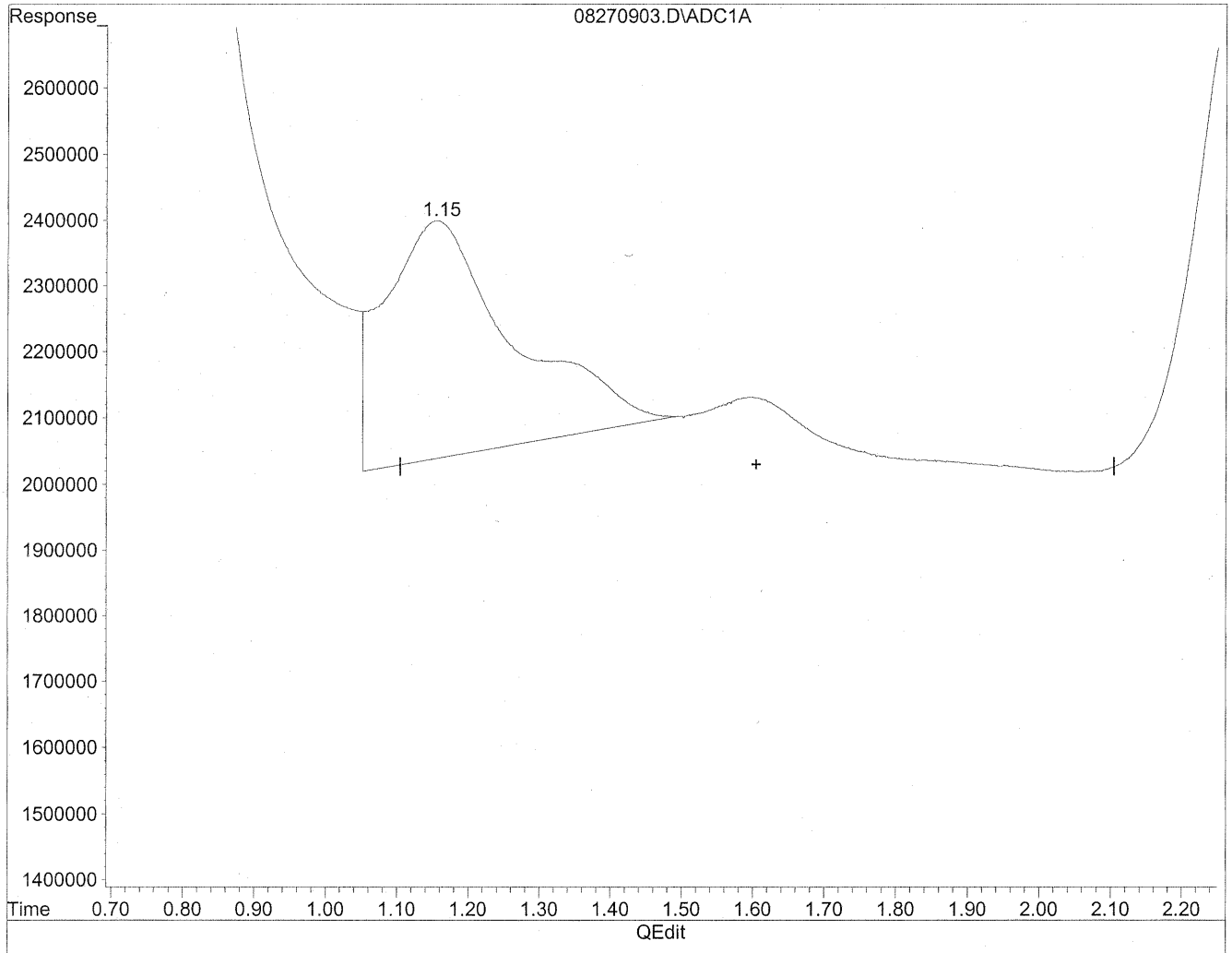
response 11166974

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270903.D Vial: 3
Acq On : 27 Aug 2009 9:35 am Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:42 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.16min 319.863ng/ml

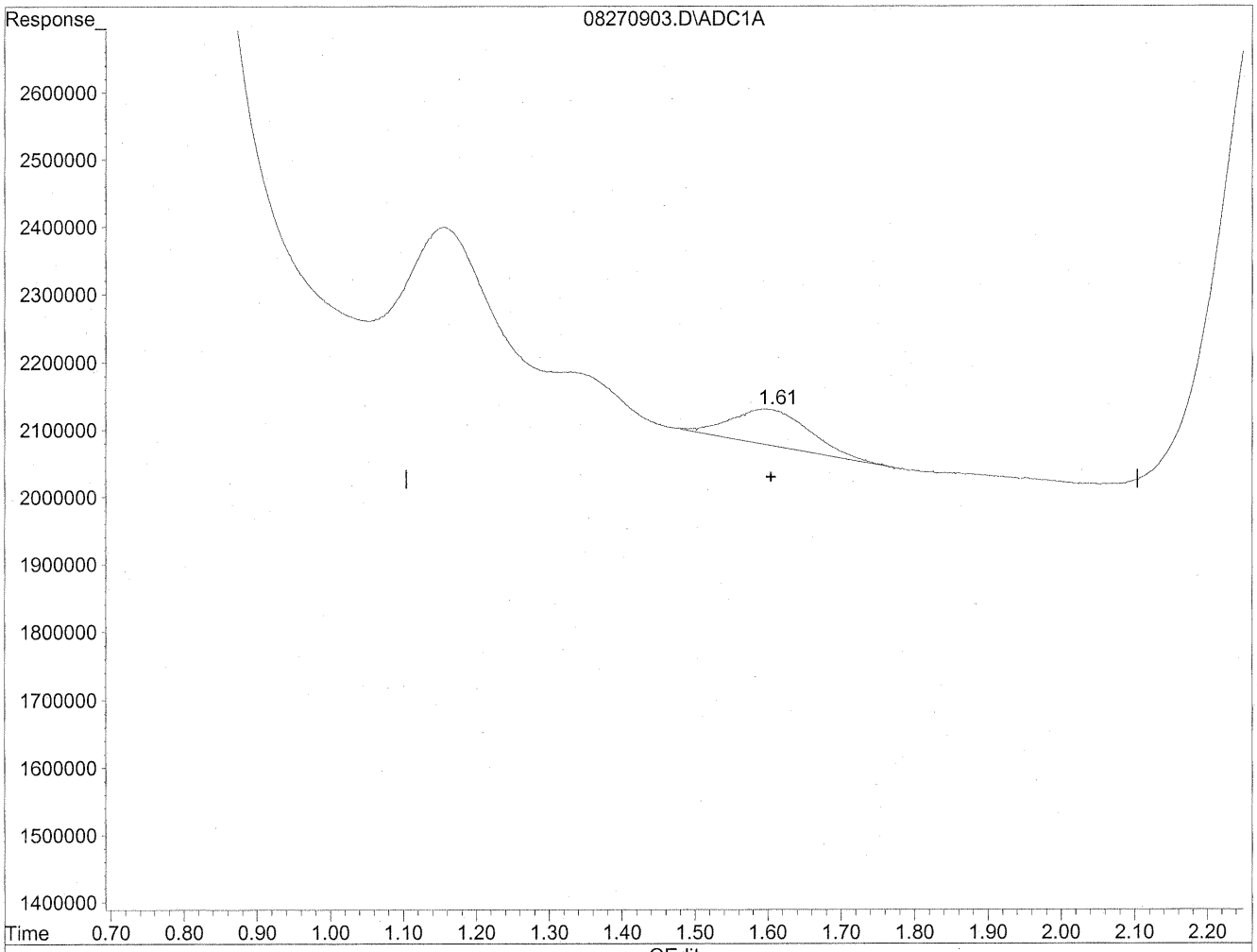
response 44852245

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270903.D Vial: 3
Acq On : 27 Aug 2009 9:35 am Operator: HC
Sample : MB front lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:42 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.61min 28.618ng/ml m

response 4012978

HL 8/31/09 LC

LM 8/31/09

182

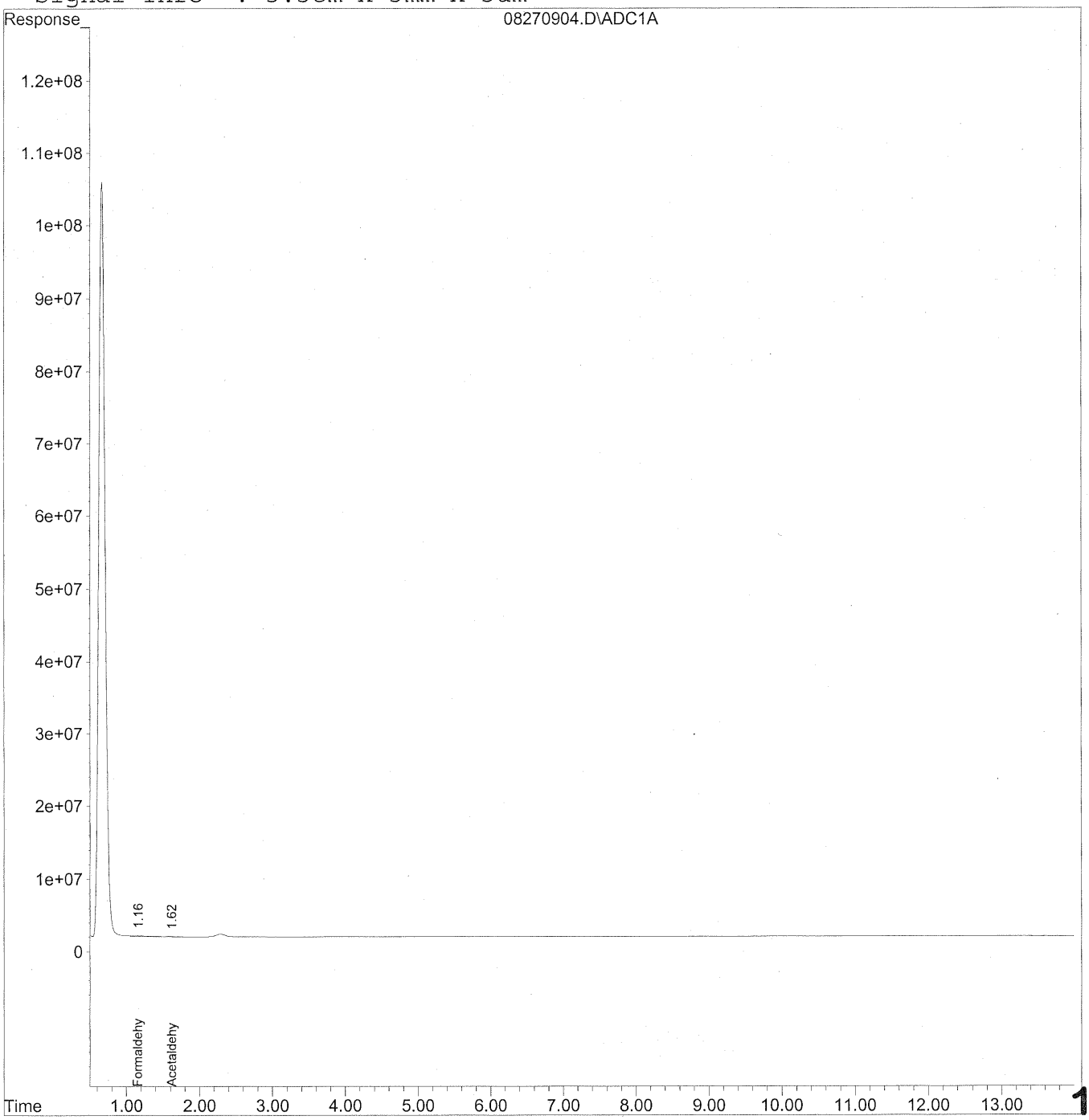
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270904.D Vial: 4
Acq On : 27 Aug 2009 9:50 am Operator: HC
Sample : MB back lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:43 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 : Sat Aug 29 16:33:38 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



183

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\27\08270904.D Vial: 4
Acq On : 27 Aug 2009 9:50 am Operator: HC
Sample : MB back lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:43 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 : Sat Aug 29 16:33:38 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

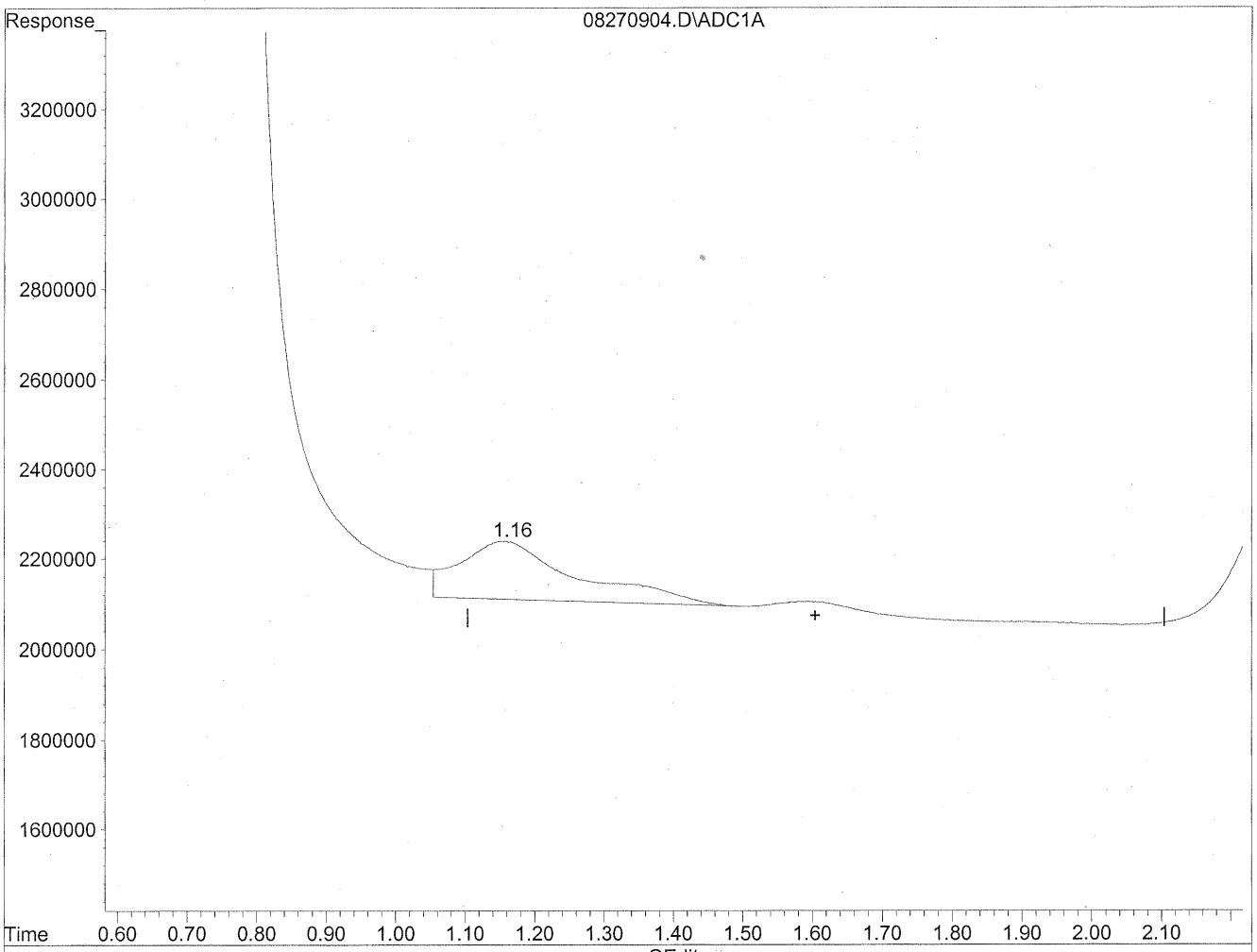
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	15097147	82.237	ng/ml
2)	Acetaldehyde	1.62	1700624	12.128	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\27\08270904.D Vial: 4
Acq On : 27 Aug 2009 9:50 am Operator: HC
Sample : MB back lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:43 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.16min 107.665ng/ml

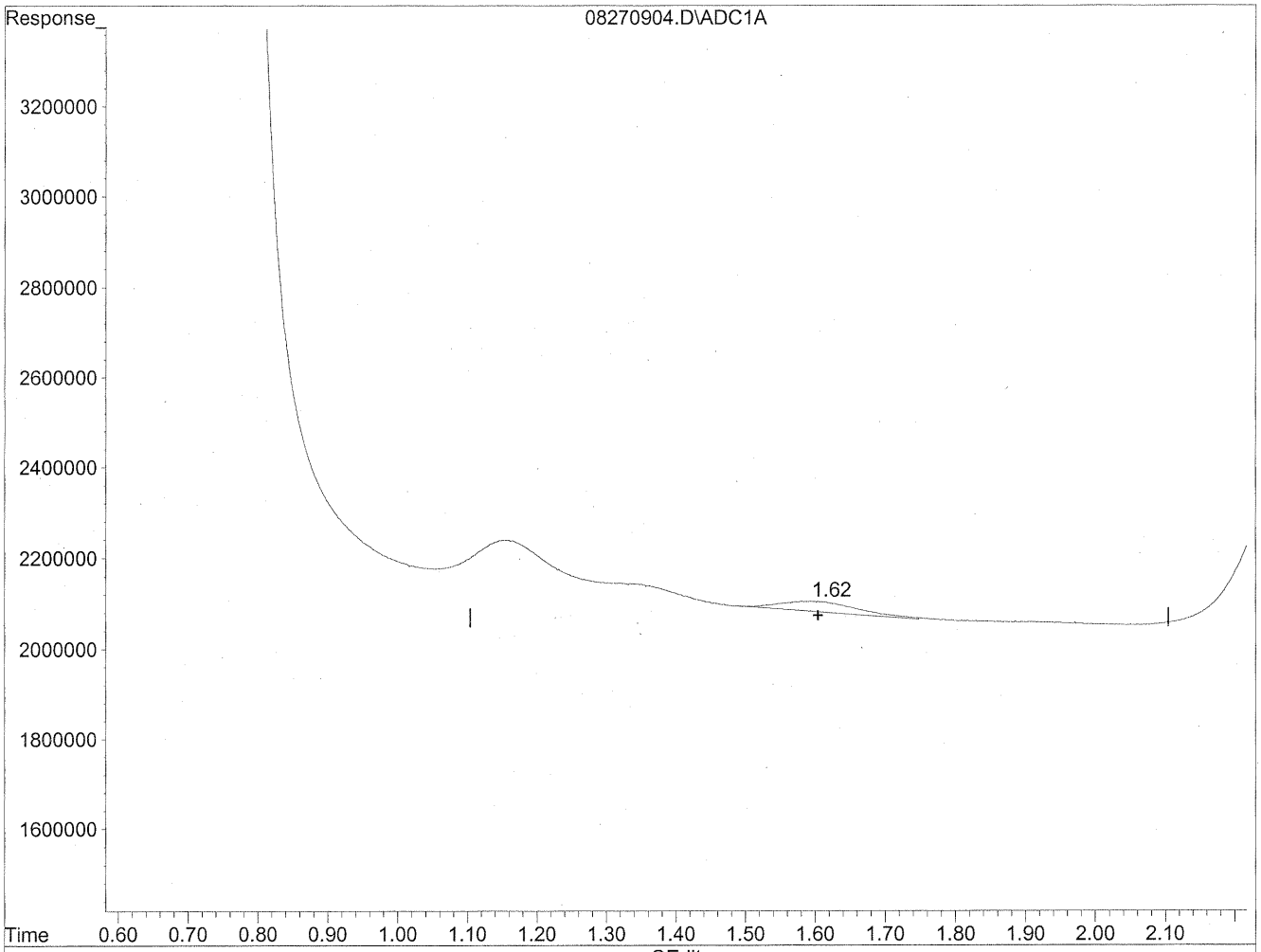
response 15097147

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270904.D Vial: 4
Acq On : 27 Aug 2009 9:50 am Operator: HC
Sample : MB back lot 5855/5994 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:43 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.62min 12.128ng/ml m

response 1700624

HC 8/31/09 WP

8/31/09

(+) = Expected Retention Time

INITIAL CALIBRATION STANDARDS

Response Factor Report LC 01

Method : J:\LC01\METHODS\TO11709B.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	4.947	4.643	4.645	4.728	4.798	4.879	E4	6.78

(#) = Out of Range

COMPOUND	50	100	500	1500	5000	10000	AVERAGE	SD	%RSD
Formaldehyde	177610.387	1.84E+05	1.82E+05	1.83E+05	1.85E+05	1.90E+05	1.84E+05	3.90E+03	2.12%
Acetaldehyde	137817.873	1.40E+05	1.40E+05	1.39E+05	1.41E+05	1.44E+05	1.40E+05	2.12E+03	1.51%
Propionaldehyde	102061.973	1.10E+05	1.06E+05	1.06E+05	1.07E+05	1.10E+05	1.07E+05	2.86E+03	2.68%
Crotonaldehyde	108243.627	9.53E+04	9.45E+04	9.44E+04	9.51E+04	9.69E+04	9.74E+04	5.38E+03	5.52%
Butyraldehyde	85497.7333	8.91E+04	8.71E+04	8.85E+04	8.91E+04	9.08E+04	8.83E+04	1.83E+03	2.07%
Benzaldehyde	61157.9267	6.91E+04	6.72E+04	6.55E+04	6.56E+04	6.67E+04	6.59E+04	2.65E+03	4.02%
Isovaleraldehyde	77804.4133	7.95E+04	7.87E+04	7.72E+04	7.76E+04	7.87E+04	7.83E+04	8.66E+02	1.11%
Valeraldehyde	76093.6467	7.70E+04	7.25E+04	7.11E+04	7.16E+04	7.28E+04	7.35E+04	2.42E+03	3.30%
o-Tolualdehyde	55101.3133	5.70E+04	5.95E+04	5.78E+04	5.97E+04	6.07E+04	5.83E+04	2.07E+03	3.55%
m,p-Tolualdehyde	50477.4933	5.57E+04	5.42E+04	5.37E+04	5.46E+04	5.54E+04	5.40E+04	1.87E+03	3.47%
Hexaldehyde	68525.6867	7.11E+04	6.46E+04	6.57E+04	6.65E+04	6.75E+04	6.73E+04	2.30E+03	3.41%
2,5-Dimethylbenzaldehyde	55134.18	4.95E+04	4.64E+04	4.65E+04	4.73E+04	4.80E+04	4.88E+04	3.31E+03	6.78%

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%	4295221	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%	9452197	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%	98878868	0.65%
1500ng/ml TO11A S	274724982	0.02%	209301649	0.12%	158919579	0.14%	142112419	0.32%	132549734	0.12%	98183657	0.06%
1500ng/ml TO11A S	273895978	0.28%	208465321	0.28%	158125683	0.36%	139629551	1.43%	131425702	0.96%	97652643	0.60%
5000ng/ml TO11A S	928364658	0.45%	70617050	0.05%	53906785	0.39%	47626854	0.19%	44639279	0.21%	32828610	0.04%
5000ng/ml TO11A S	925768000	0.17%	70855241	0.38%	54013392	0.59%	47784449	0.52%	44656805	0.25%	32841355	0.08%
5000ng/ml TO11A S	918424042	0.62%	70279188	0.43%	53167508	0.98%	47195457	0.72%	44344183	0.45%	32776290	0.12%
10000ng/ml TO11A	1908653125	0.62%	1450154617	0.67%	1099941045	0.36%	972691462	0.37%	91089701	0.36%	66846212	0.28%
10000ng/ml TO11A	1905913073	0.48%	1446499891	0.41%	1098837646	0.26%	971357788	0.23%	91132824	0.41%	66912896	0.38%
10000ng/ml TO11A	1875917434	1.10%	1425028469	1.08%	1089338811	0.61%	963283335	0.60%	90056123	0.78%	66223844	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%	3387183	22.94%	5445142	7.87%	3244418	5.31%	2546144	7.64%
50ng/ml TO11A Std	4002738	2.89%	4025564	5.81%	2461625	10.65%	4897087	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%	10979457	1.36%	6702769	5.76%	5399082	9.13%
100ng/ml TO11A St	8338385	4.88%	8117341	5.49%	5921917	3.82%	11235135	0.94%	7714022	8.46%	4735227	4.29%
100ng/ml TO11A St	8025579	0.95%	7906862	2.75%	5642221	1.09%	11177259	0.42%	6920120	2.70%	4707951	4.84%
500ng/ml TO11A St	37944016	3.60%	35574509	1.84%	29317615	1.49%	53274975	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%	68873541	1.15%
1500ng/ml TO11A S	114690000	0.92%	105937177	0.73%	87824227	1.30%	159292531	1.13%	98846718	0.24%	70224595	0.79%
5000ng/ml TO11A S	388247386	0.05%	357832844	0.04%	298513860	0.05%	545640330	0.02%	332315493	0.11%	235692401	0.30%
5000ng/ml TO11A S	388941560	0.23%	359676615	0.47%	300077384	0.48%	547211501	0.27%	333701808	0.31%	237108293	0.30%
5000ng/ml TO11A S	386992833	0.28%	356464469	0.43%	297374461	0.43%	544331756	0.26%	332038452	0.19%	236428207	0.01%
10000ng/ml TO11A	790328317	0.44%	730218673	0.36%	608208276	0.16%	1111180147	0.26%	673516807	0.25%	478460947	0.27%
10000ng/ml TO11A	788026190	0.15%	729839210	0.31%	610326238	0.50%	1113209810	0.45%	681915785	0.99%	484763918	1.04%
10000ng/ml TO11A	782256804	0.59%	722749626	0.67%	603256599	0.66%	1100384573	0.71%	670193360	0.74%	476113656	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	4947420
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

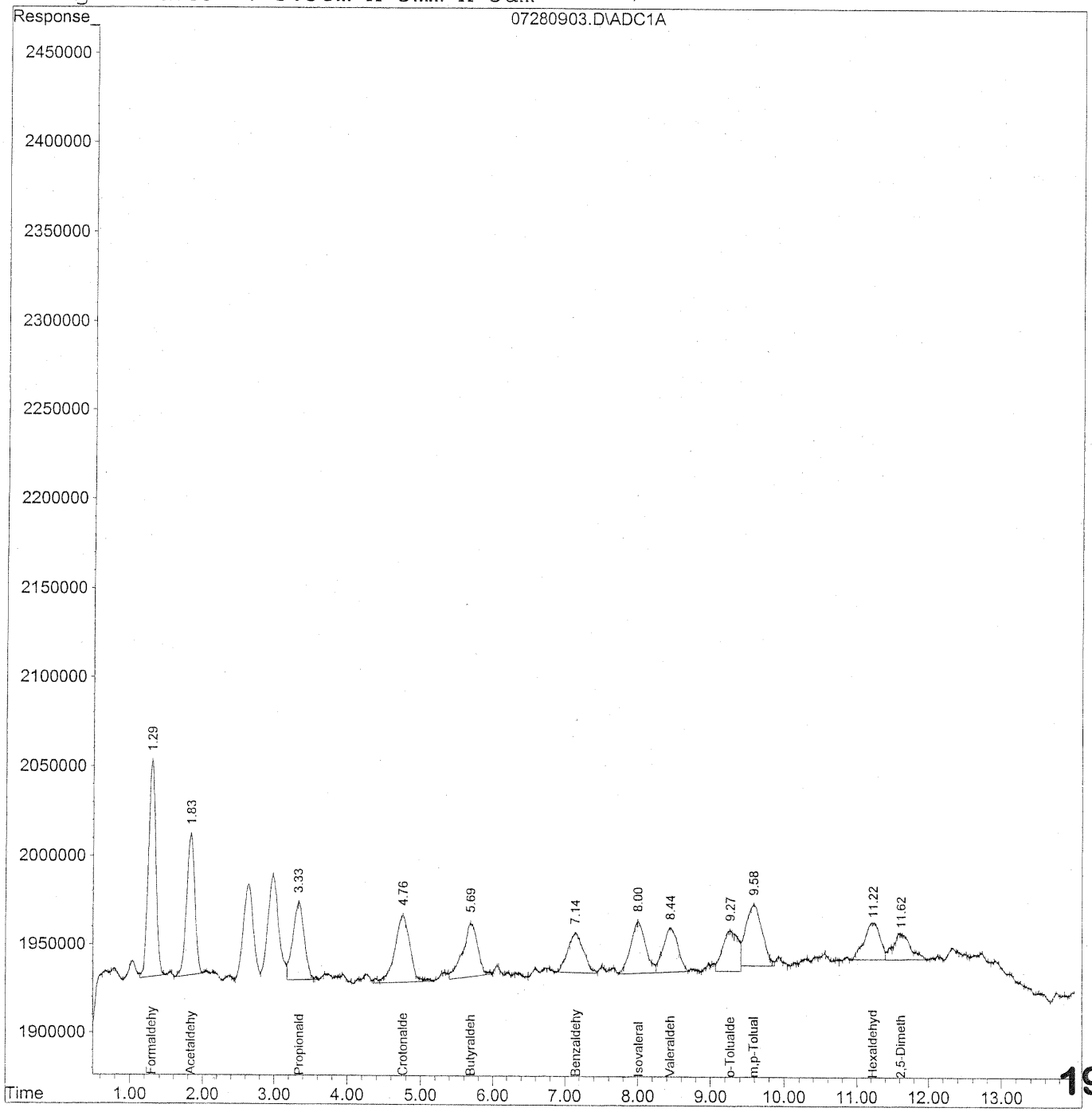
TO-11A CALIBRATION STANDARDS LIST							
50ng/ml TO11A Std S21-07270908							
100ng/ml TO11A Std S21-07270905							
500ng/ml TO11A Std S21-07270904							
1500ng/ml TO11A Std S21-07270903							
5000ng/ml TO11A Std S21-07270902							
10000ng/ml TO11A Std S21-07270901							

HC
7/29/05

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



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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

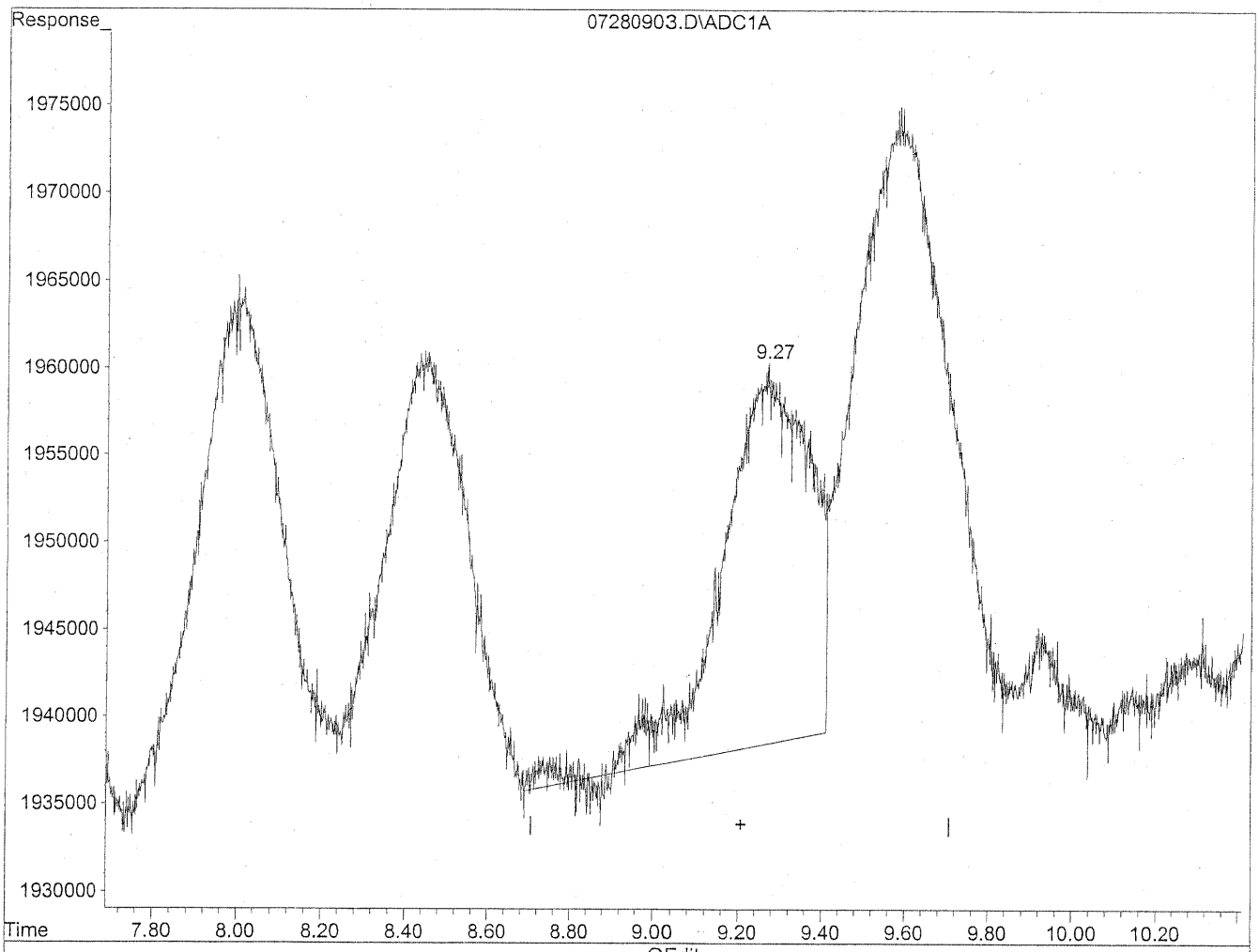
	Compound	R.T.	Response	Conc Units
	Target Compounds			
1)	Formaldehyde	1.29	8477013	48.277 ng/ml
2)	Acetaldehyde	1.83	6307171	46.755 ng/ml
3)	Propionaldehyde	3.34	4892636	47.596 ng/ml
4)	Crotonaldehyde	4.76	5507079	49.813 ng/ml
5)	Butyraldehyde	5.70	4412295	54.828 ng/ml
6)	Benzaldehyde	7.15	3362429	53.310 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

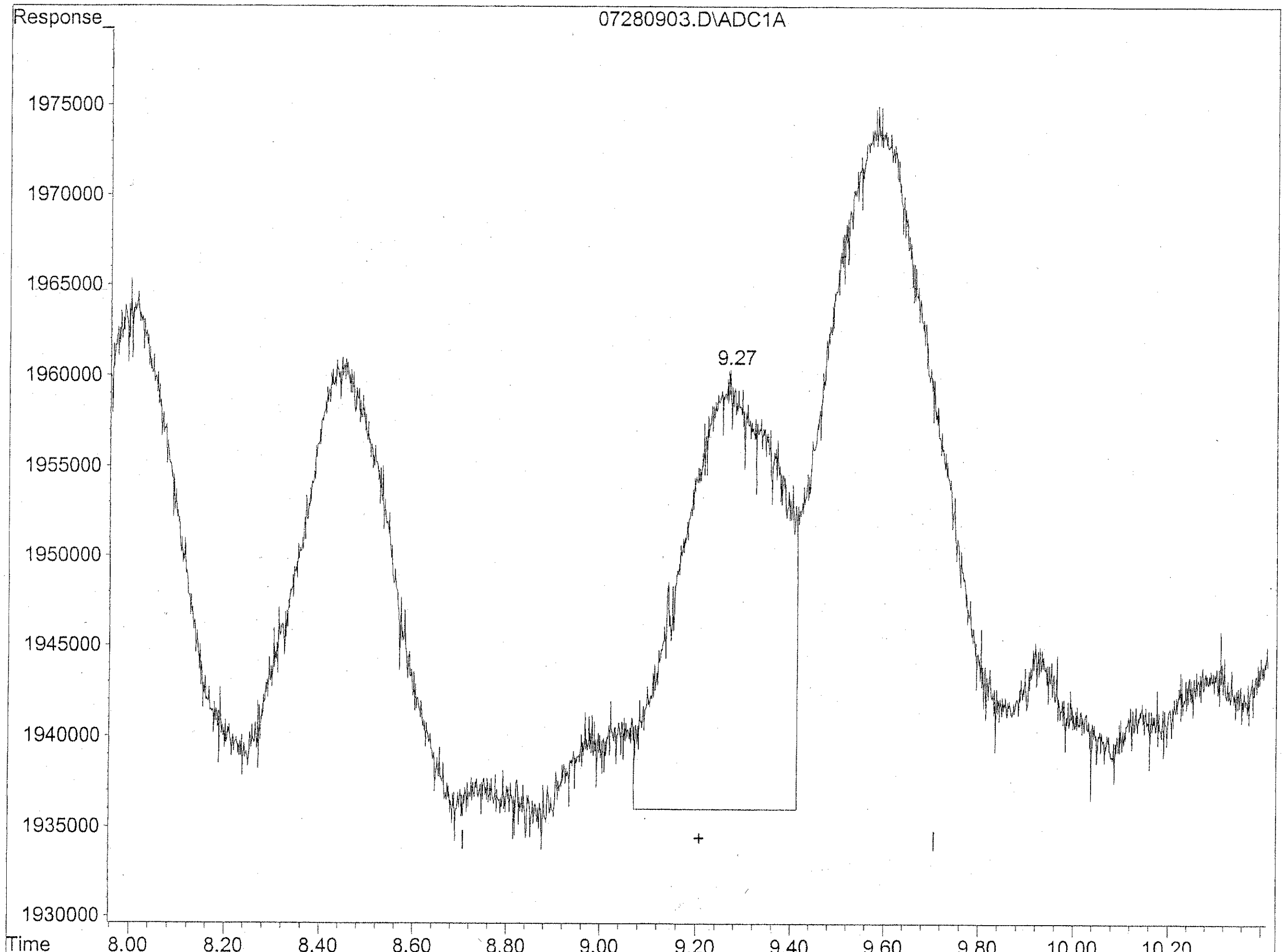
198

(+) = 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
```



QEdit

(9) o-Tolualdehyde

9.27min 62.877ng/ml m

response 3387183

HC 7/28/09 IC

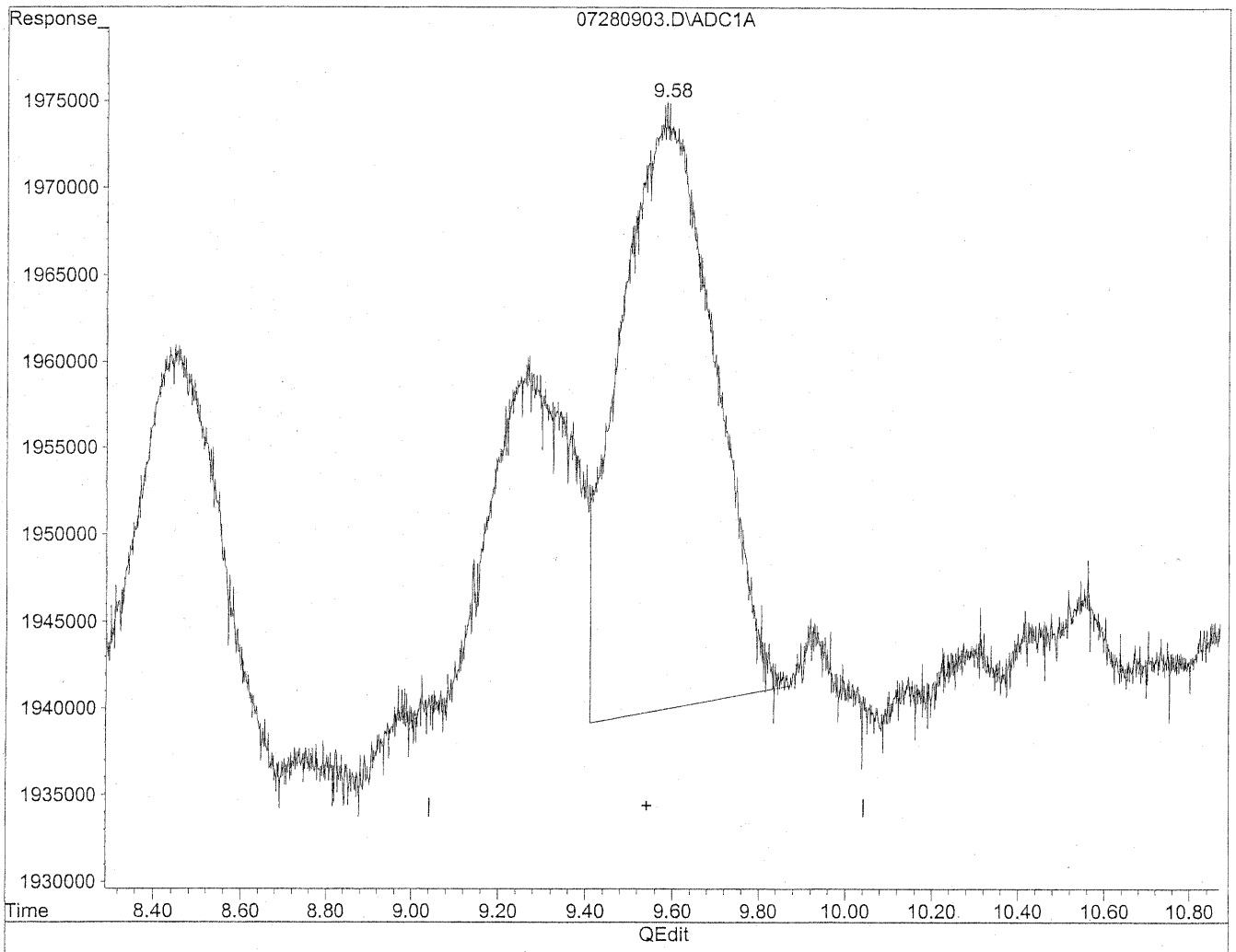
KR 7/29/09

(+) = 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

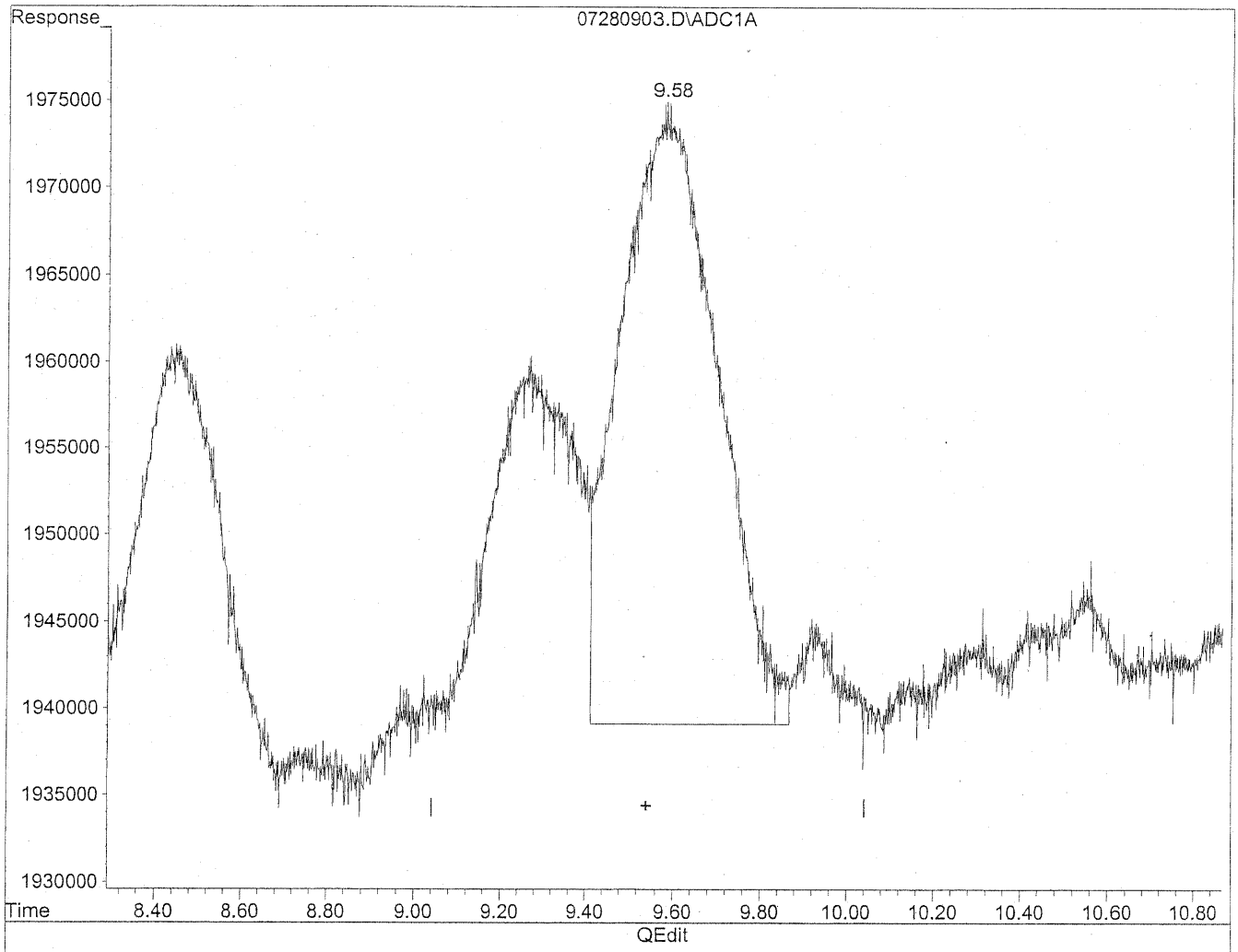
200

(+) = 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

HC 7/28/09 BC

7/29/09

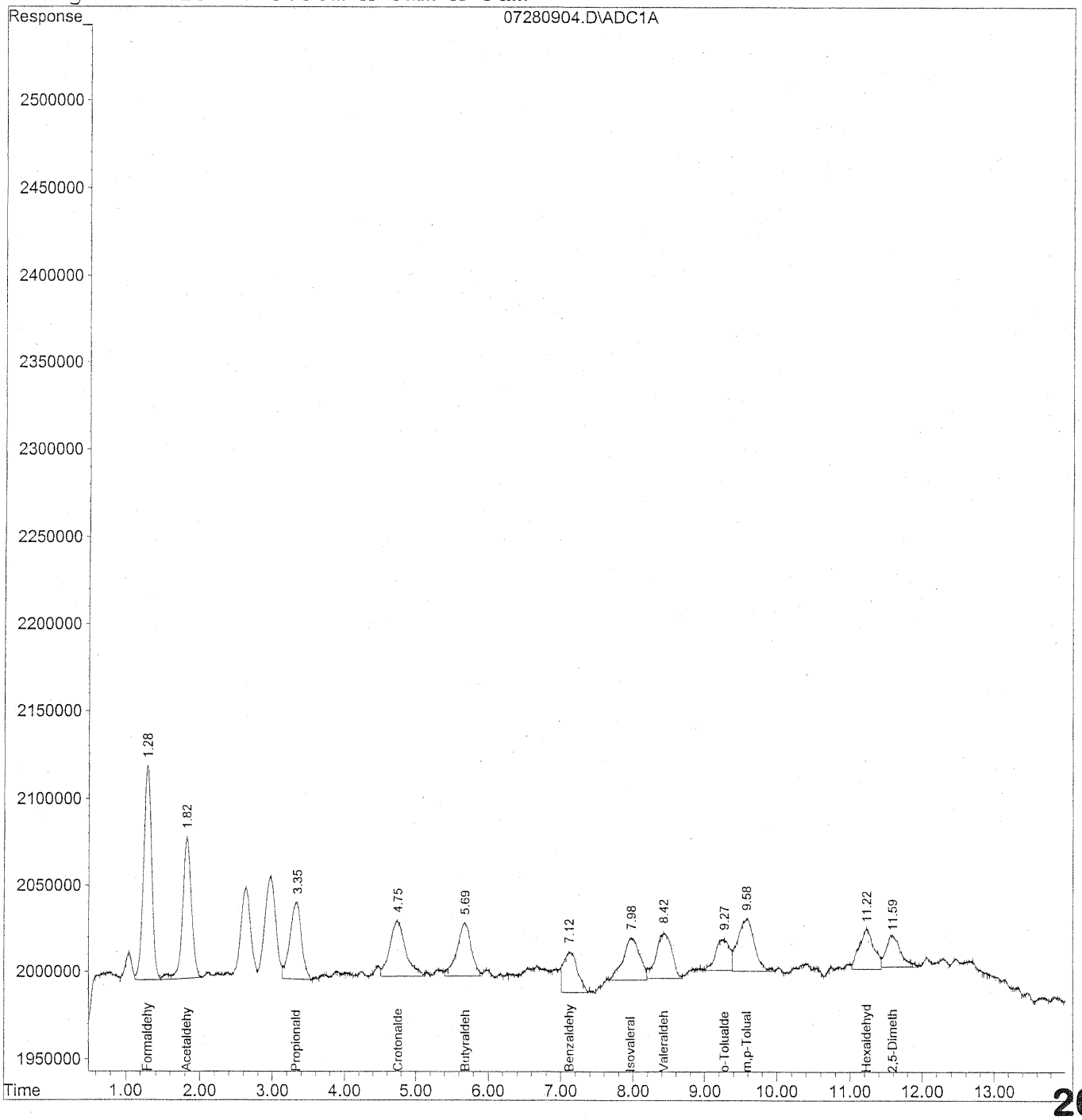
201

(+) = 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

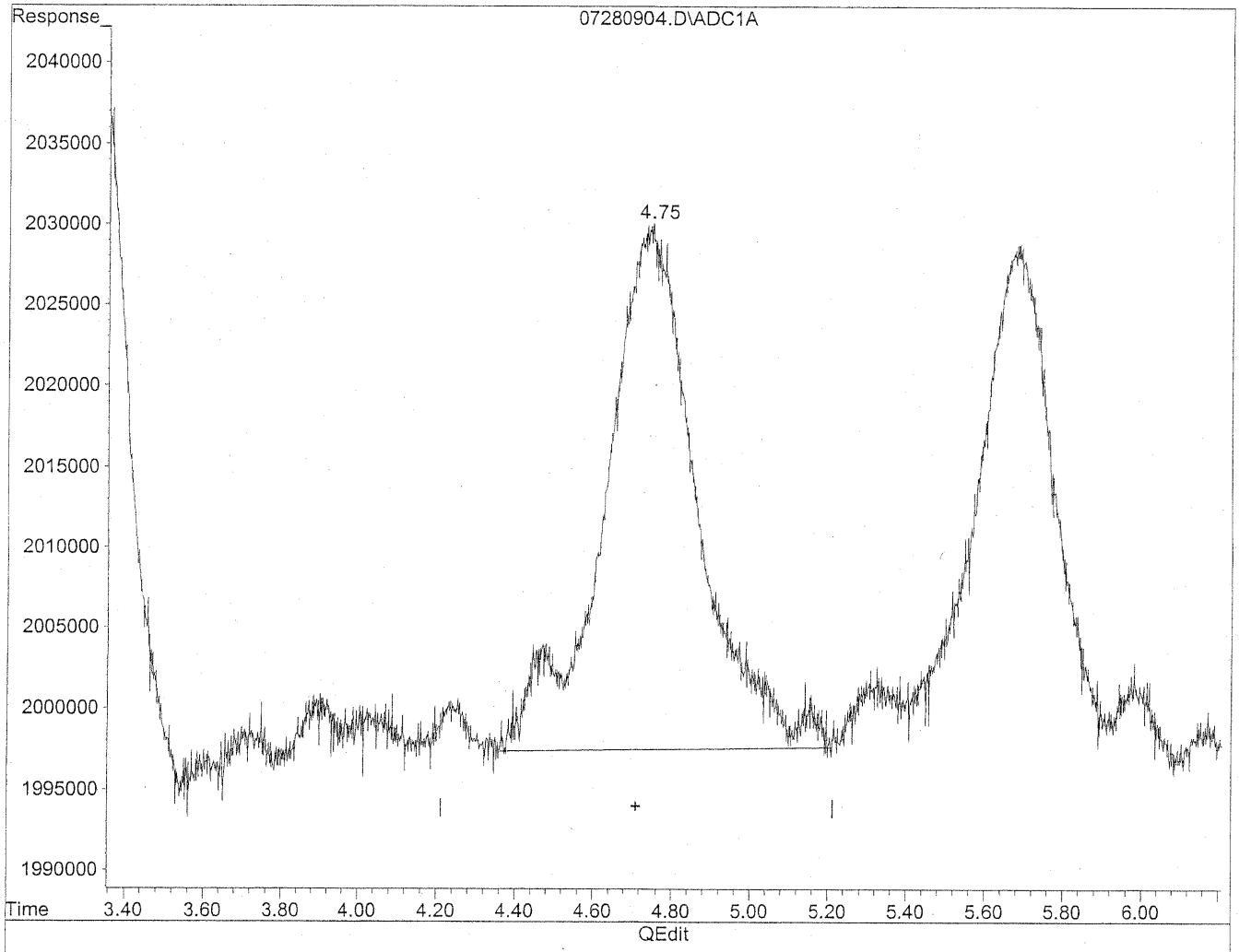
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	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

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

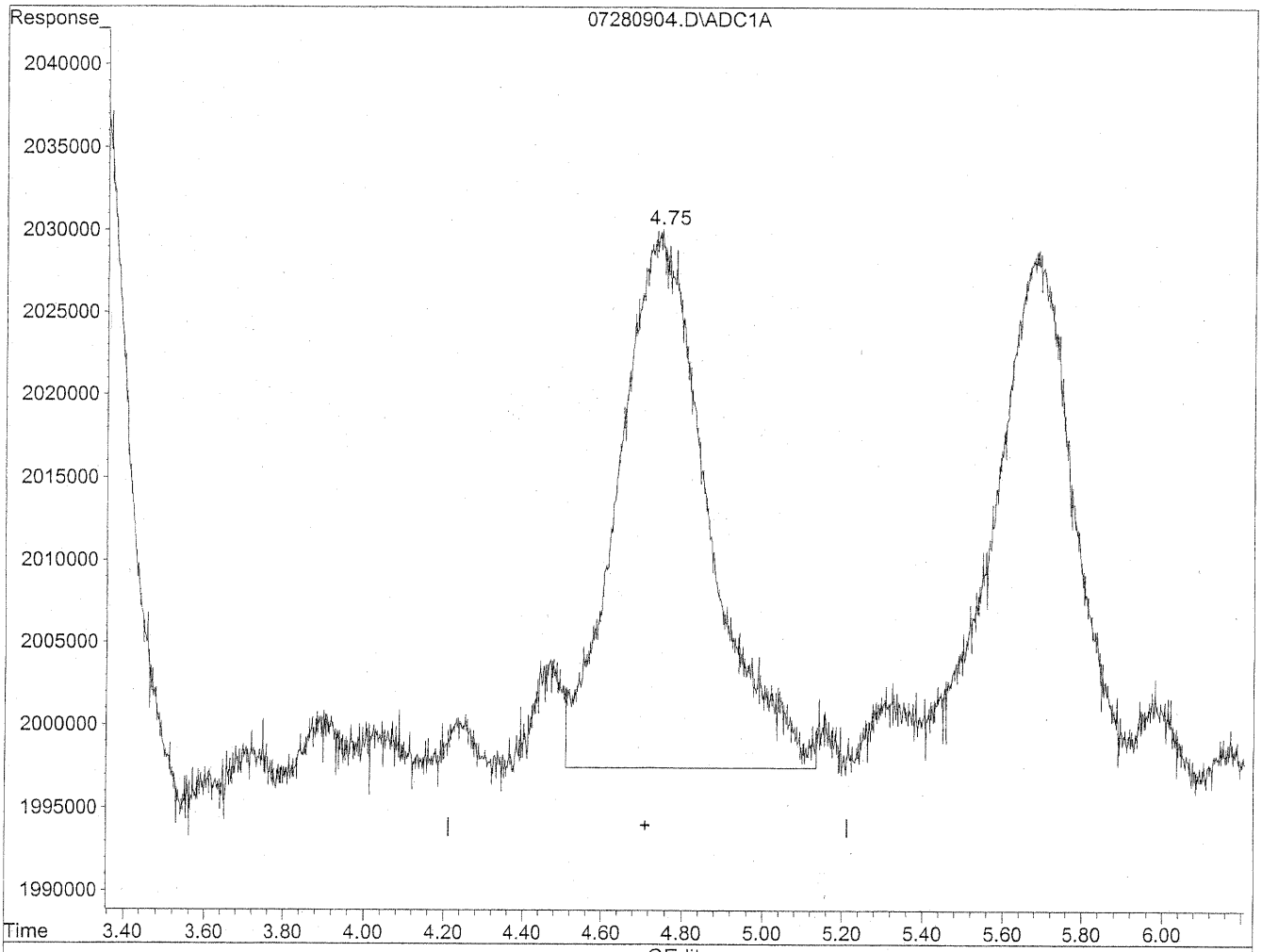
204

(+) = 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

205

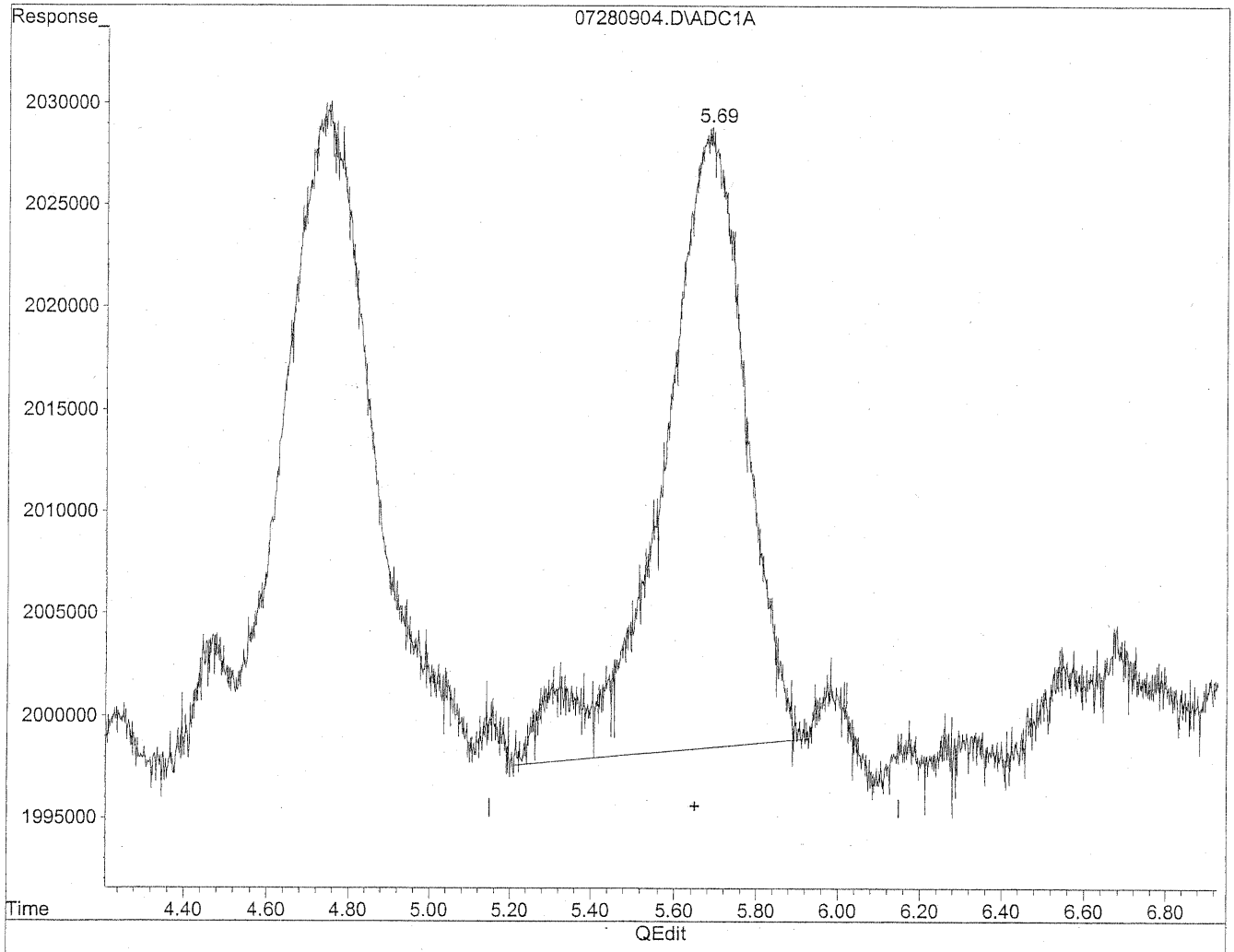
(+) = Expected Retention Time

07280904.D TO110709.M Tue Jul 28 14:23:40 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

5.68min 53.153ng/ml

response 4277470

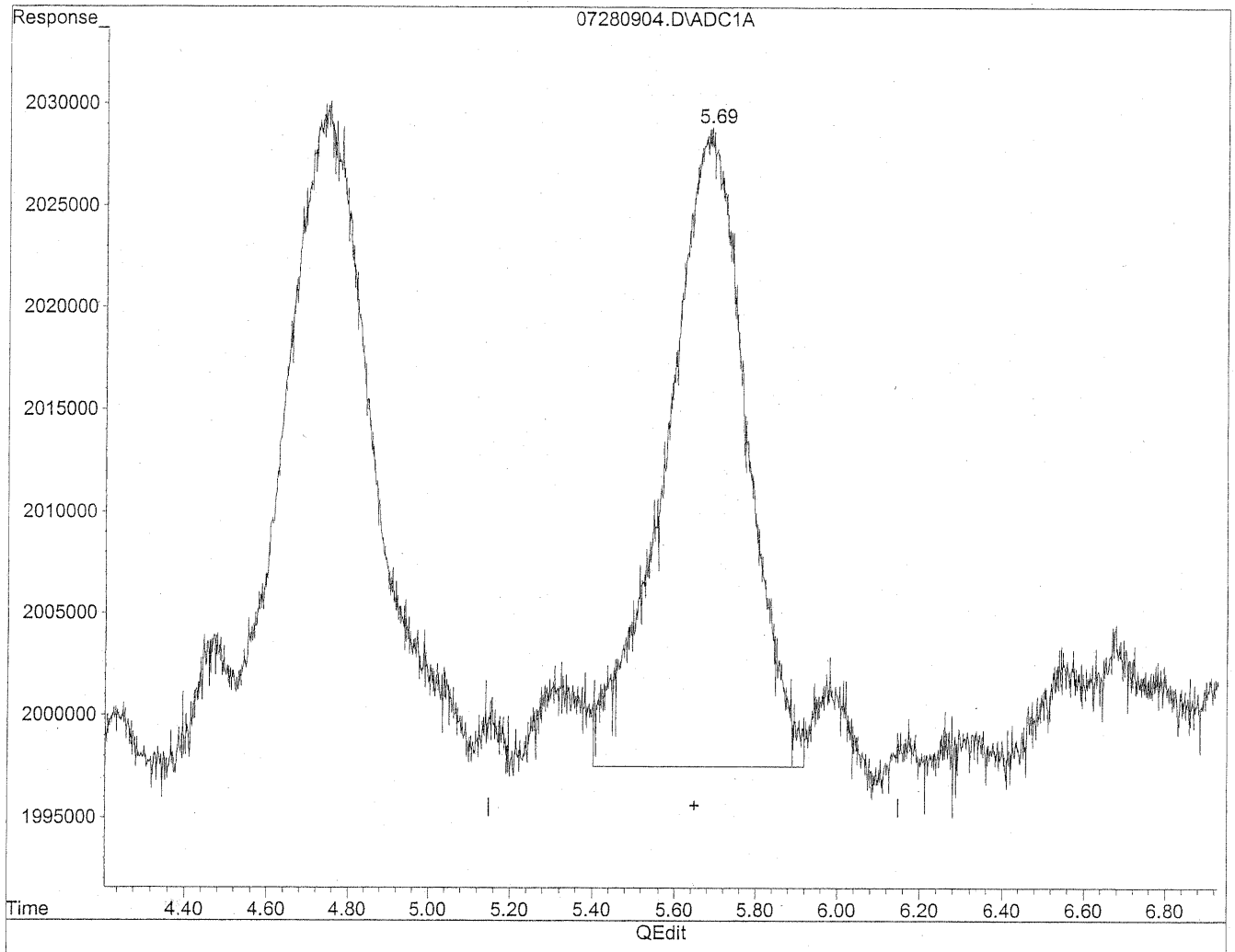
(+) = Expected Retention Time

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 7/28/09
SH

KR 7/29/09

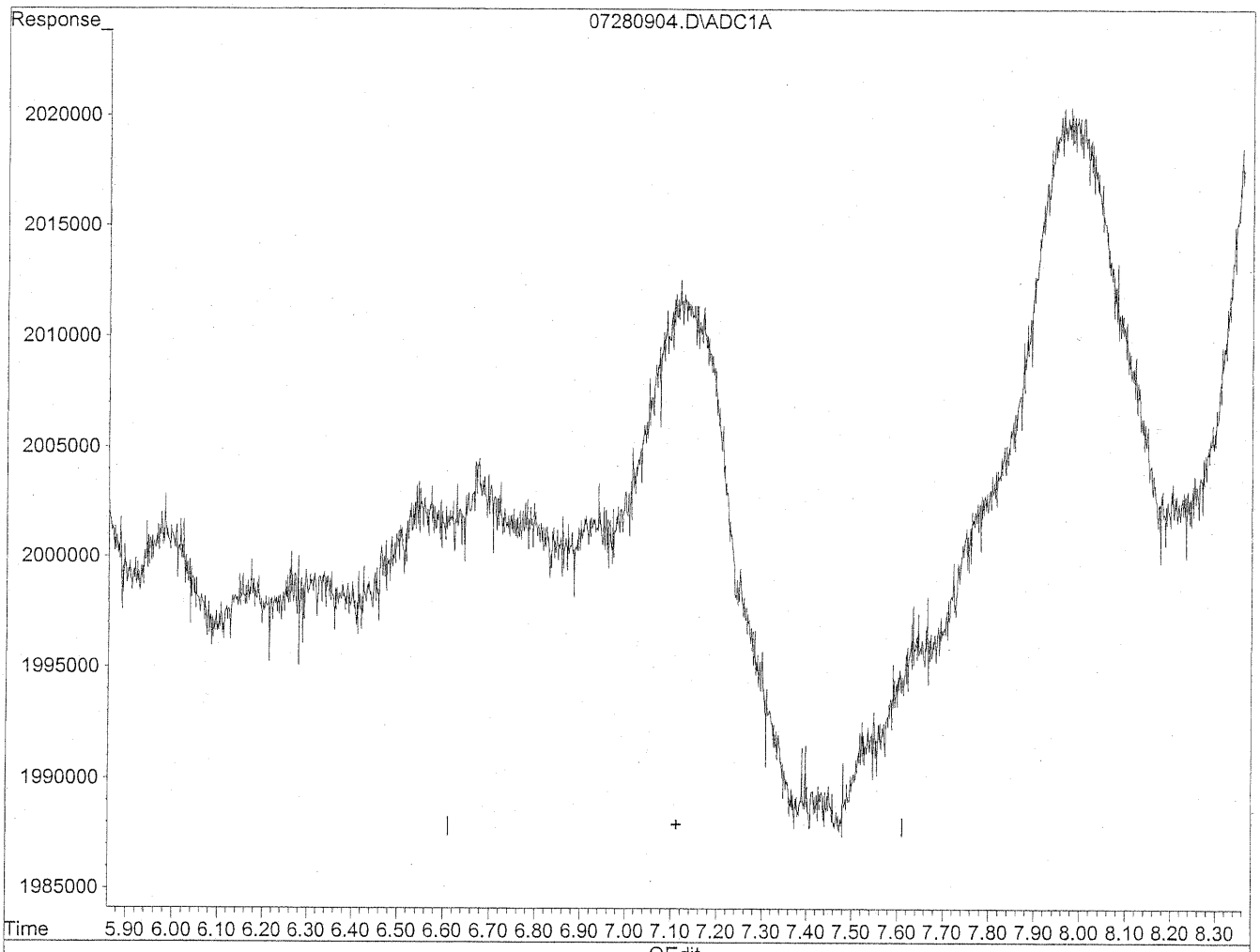
207

(+) = Expected Retention Time
07280904.D TO110709.M Tue Jul 28 14:23:54 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

(6) Benzaldehyde

7.11min 0.000ng/ml

response 0

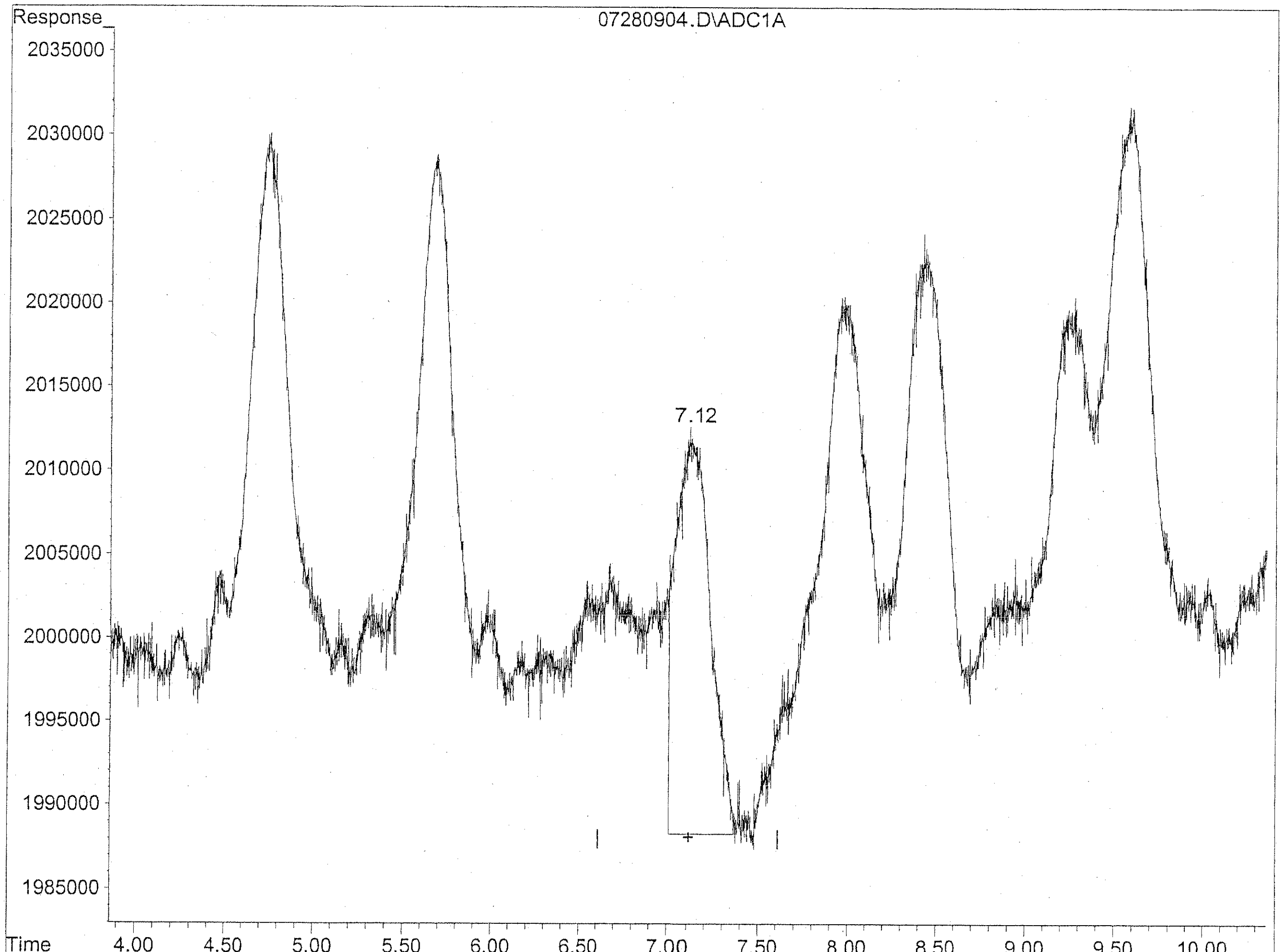
208

(+) = 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



QEdit

(6) Benzaldehyde

7.12min 48.820ng/ml m

response 3079204

HC 7/28/09 BN1

KR 7/29/09

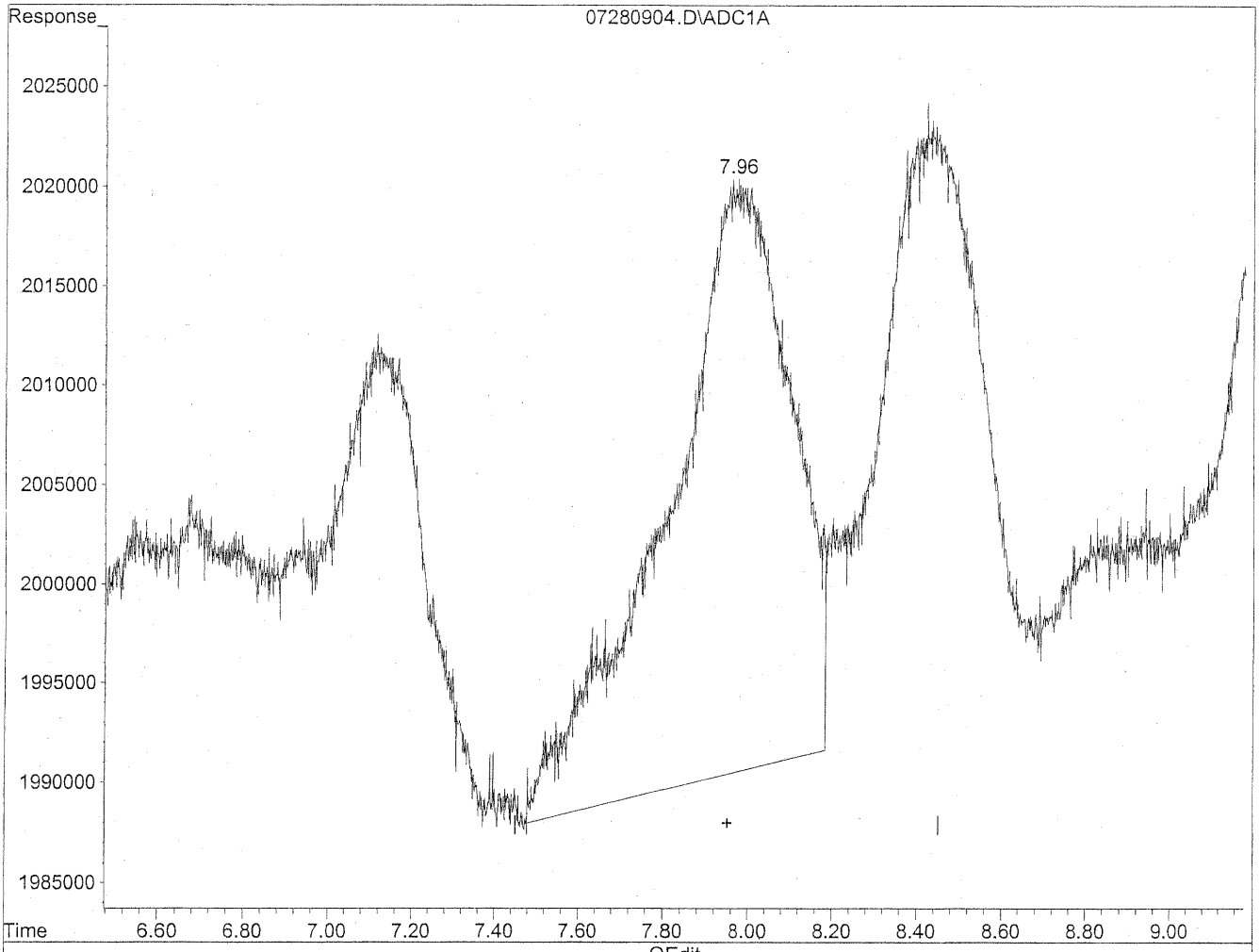
(+) = Expected Retention Time

07280904.D TO110709.M Tue Jul 28 14:24:15 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



(7) Isovaleraldehyde

7.97min 68.251ng/ml

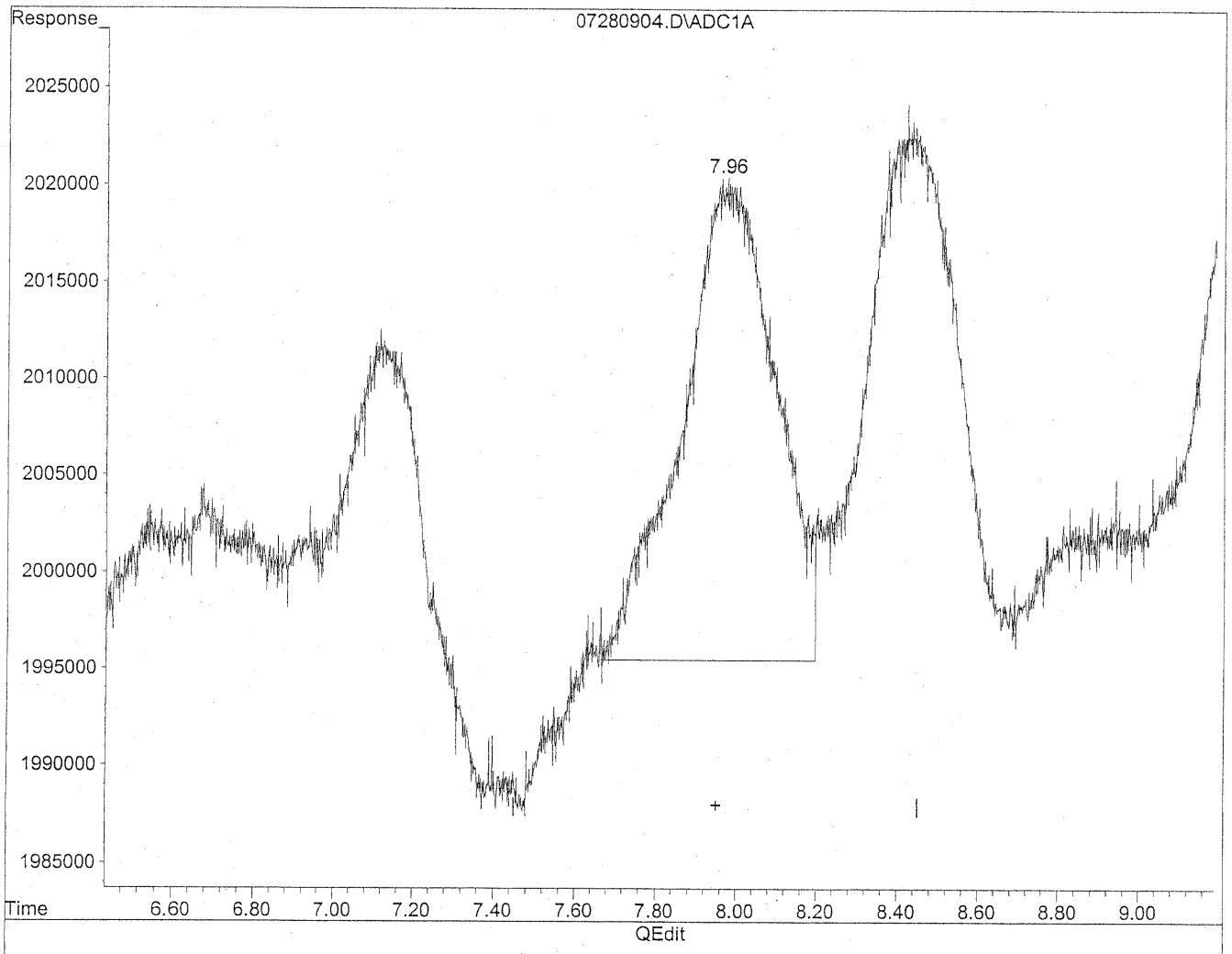
response 6050534

(+) = 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



(7) Isovaleraldehyde

7.96min 45.151ng/ml m

response 4002738

KC 7/29/09 LC

KC 7/30/09

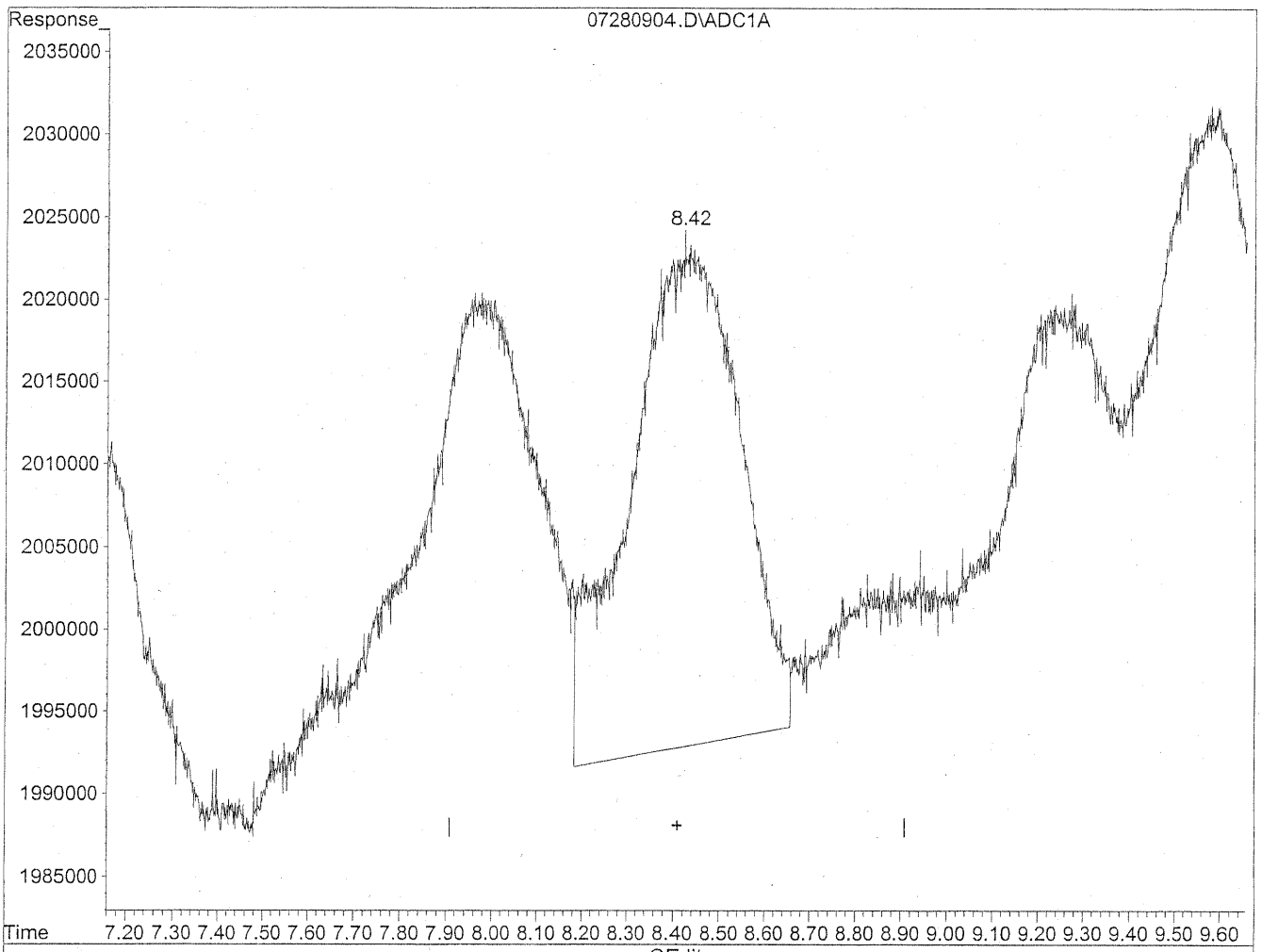
211

(+) = 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



(8) Valeraldehyde

8.43min 61.279ng/ml

response 5091976

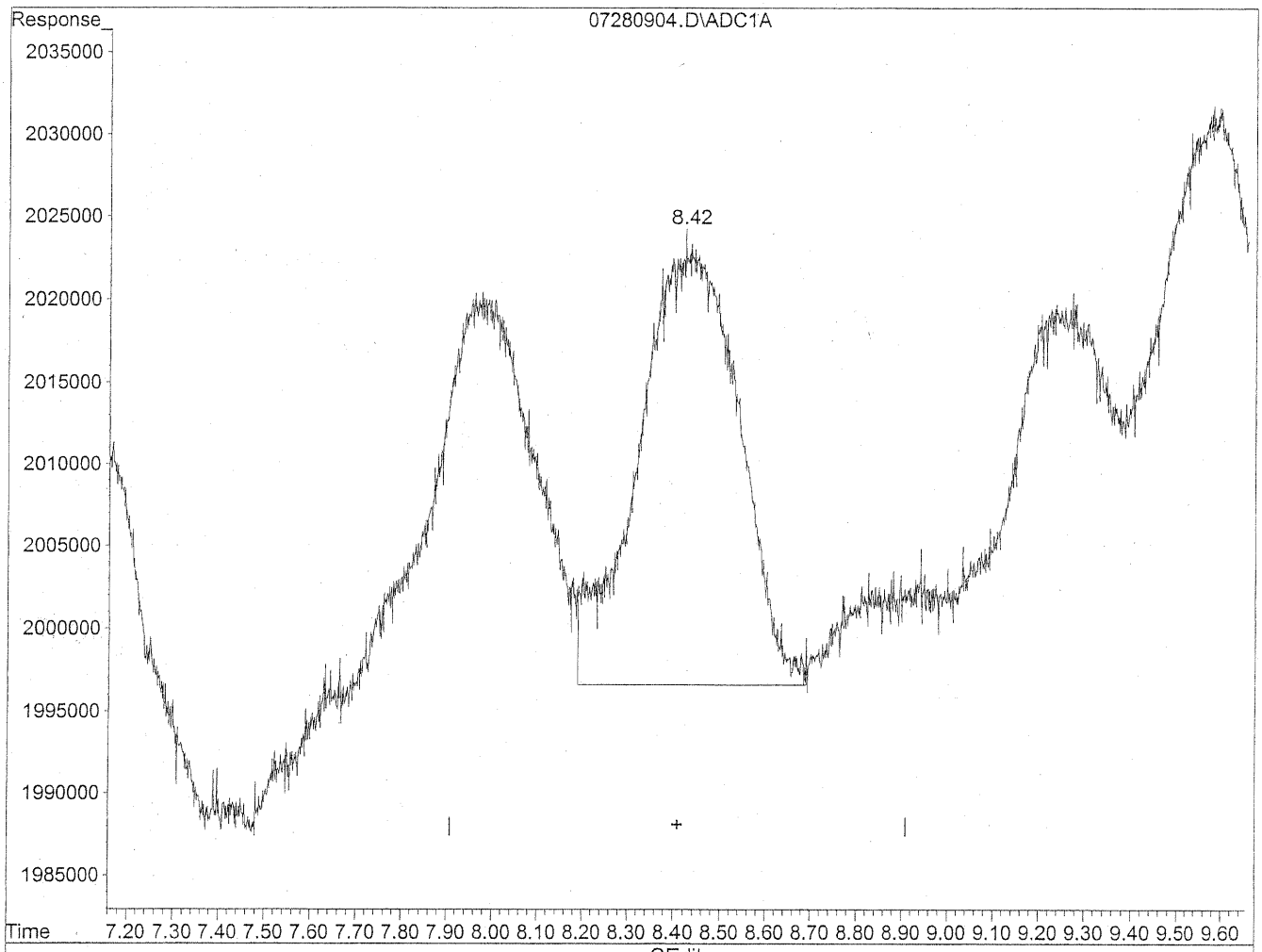
212

(+) = 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



QEdit

(8) Valeraldehyde

8.42min 48.445ng/ml m

response 4025564

HC station 1C

KR 7/29/09

213

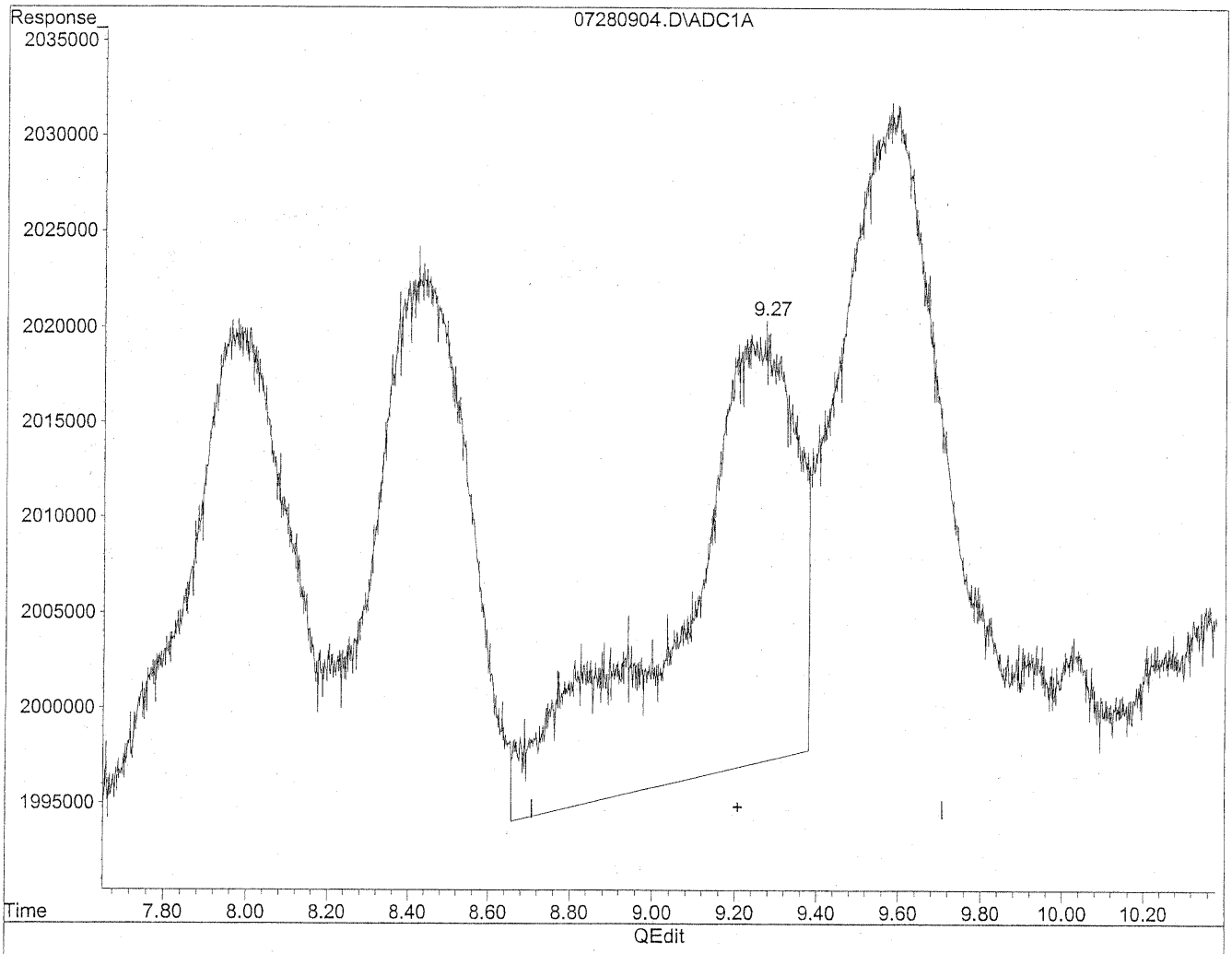
(+) = Expected Retention Time

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

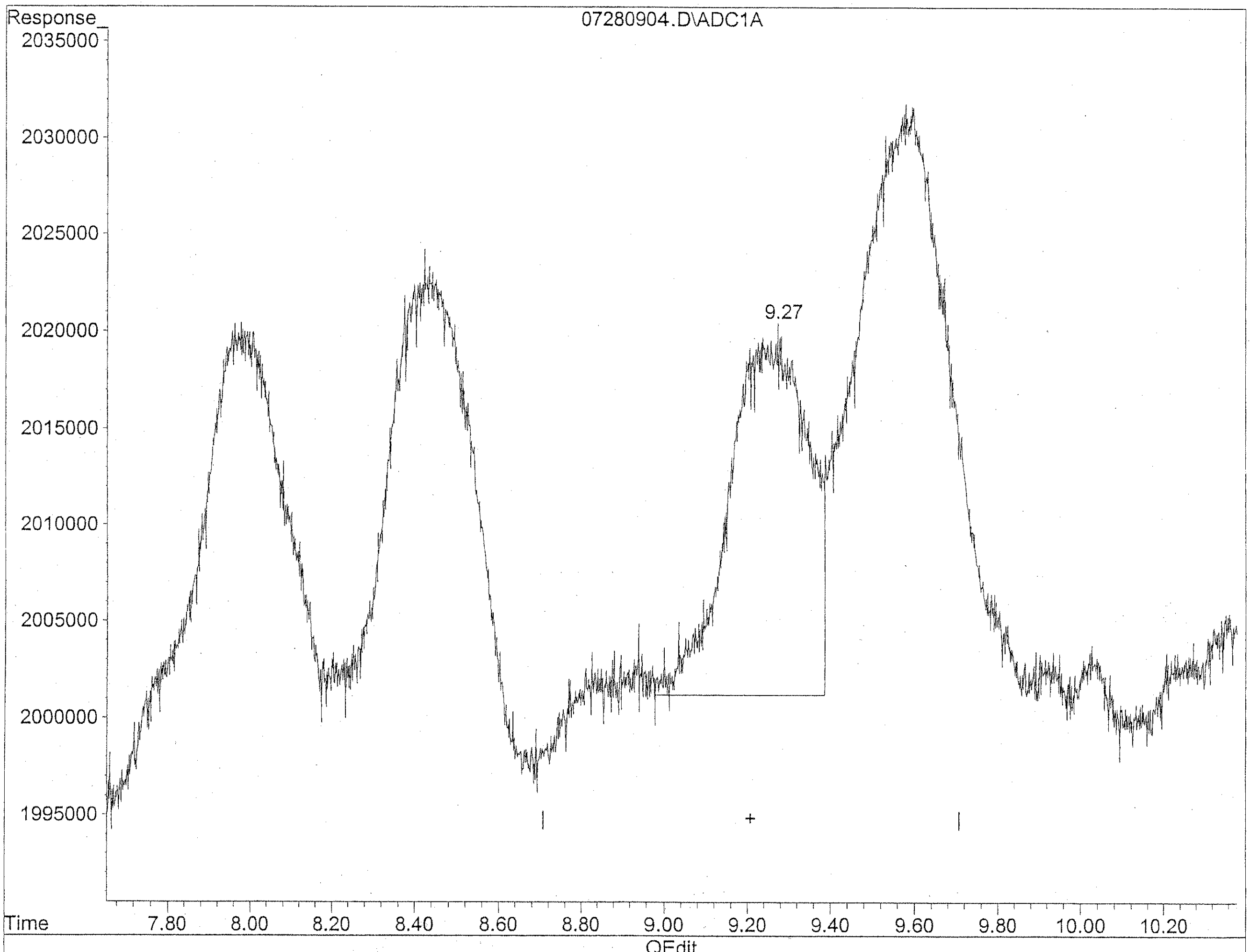
214

(+) = 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

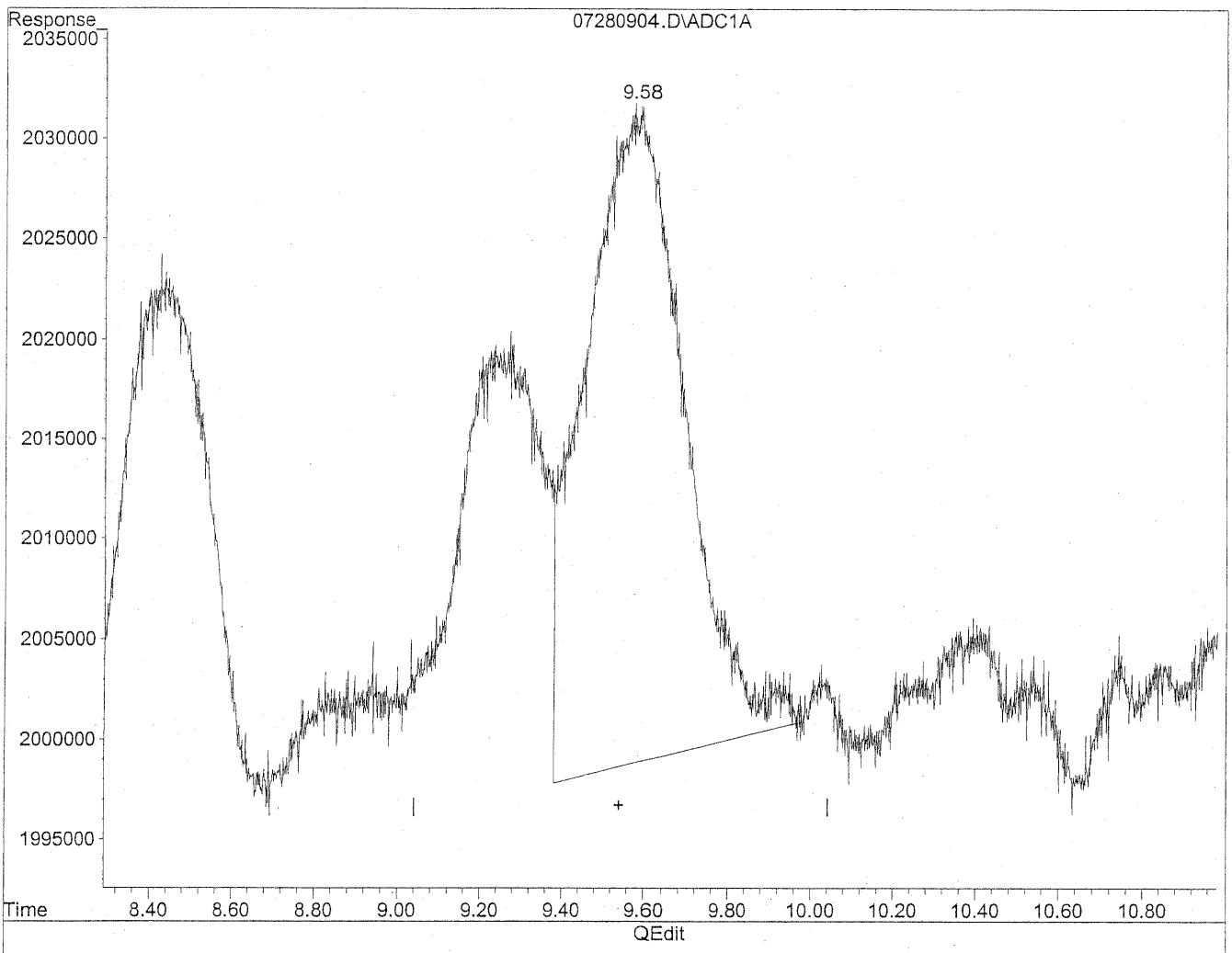
(+) = Expected Retention Time

07280904.D   TO110709.M   Tue Jul 28 14:24:57 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



(10) m,p-Tolualdehyde

9.59min 100.987ng/ml

response 5439618

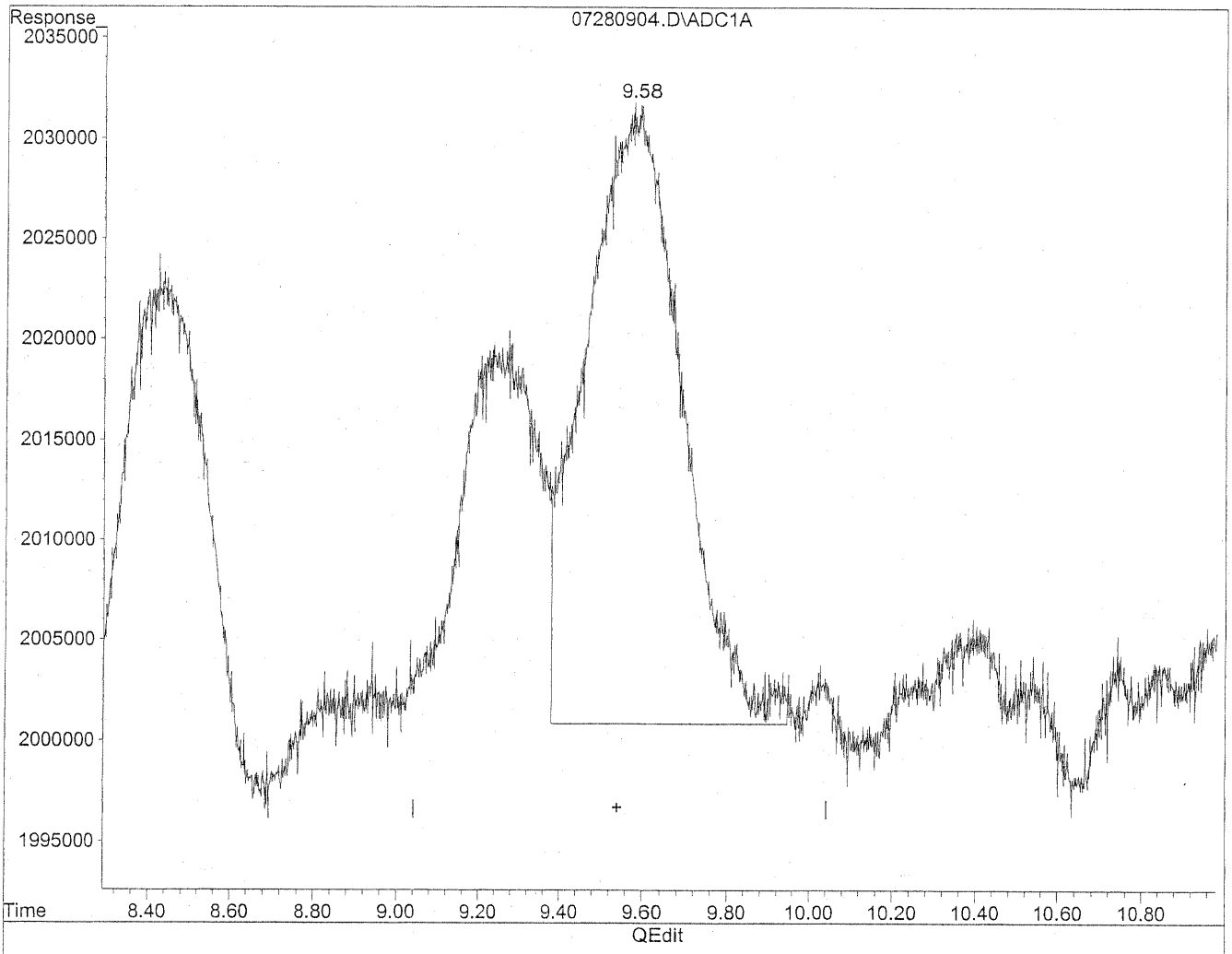
216

(+) = 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

217

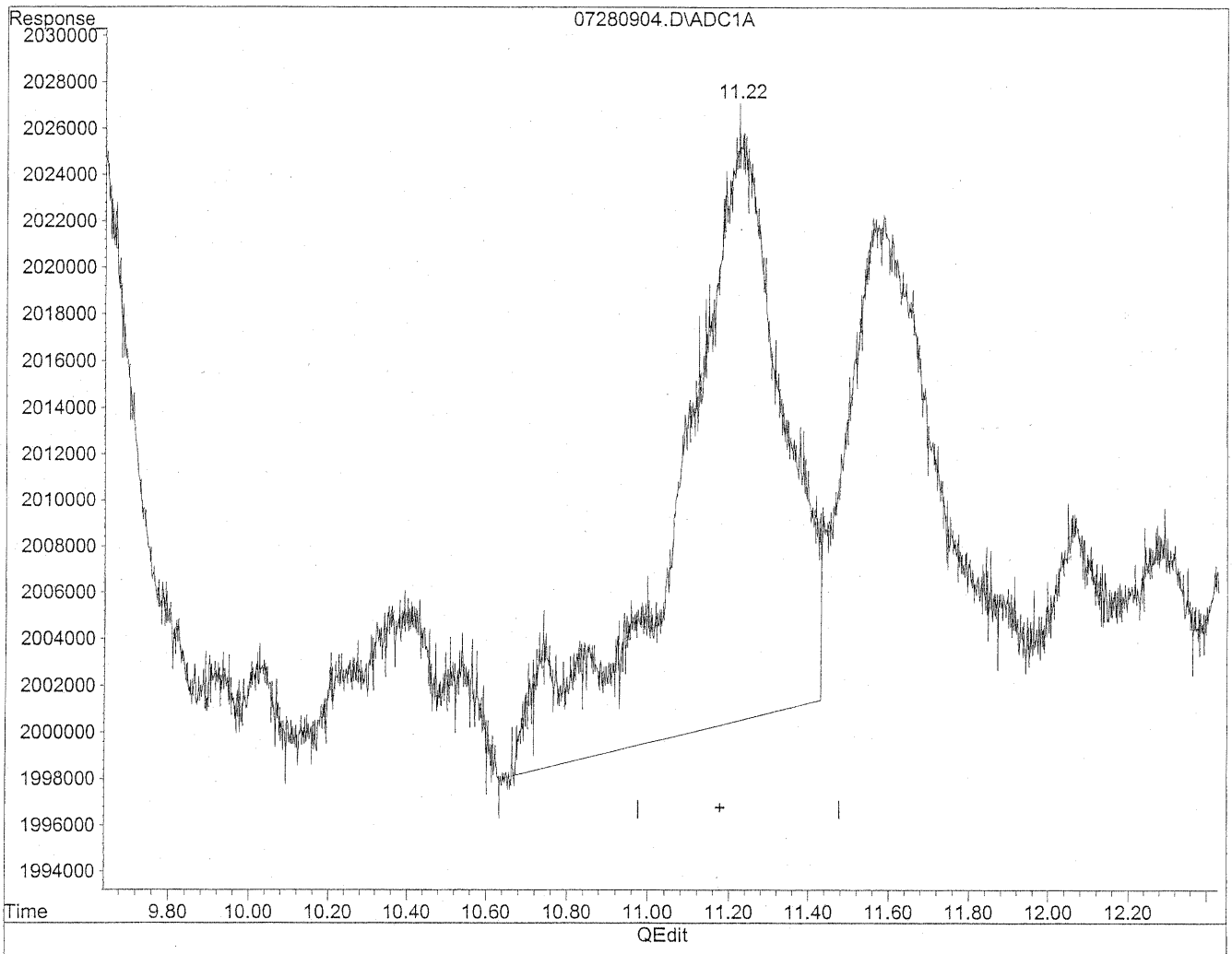
(+) = Expected Retention Time

07280904.D TO110709.M Tue Jul 28 14:25:08 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



(11) Hexaldehyde

11.23min 66.912ng/ml

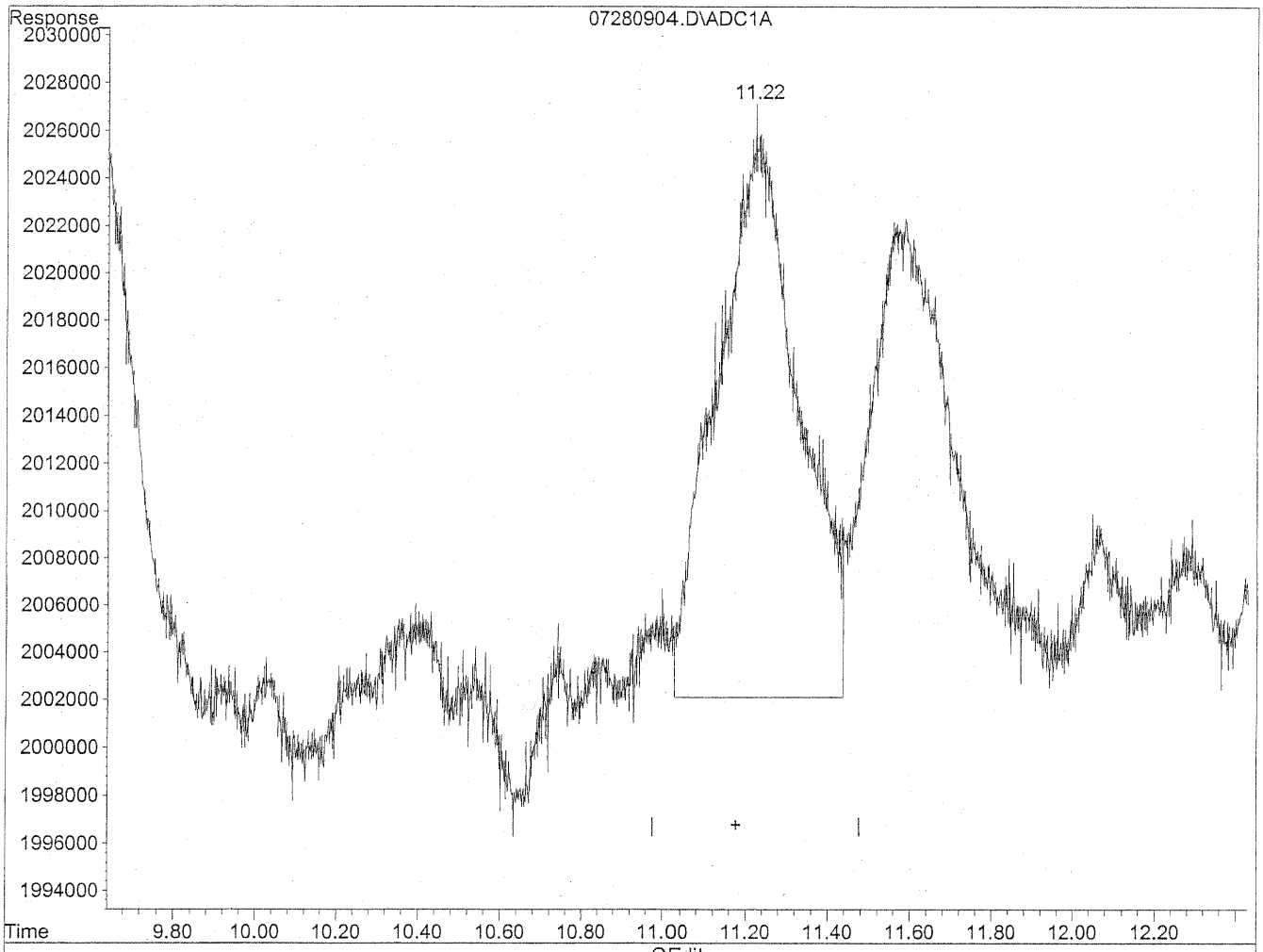
response 4492347

(+) = Expected Retention Time
07280904.D TO110709.M Tue Jul 28 14:25:12 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

(11) Hexaldehyde

11.22min 49.079ng/ml m

response 3295067

HC 7/28/09 SH

KL 7/29/09

219

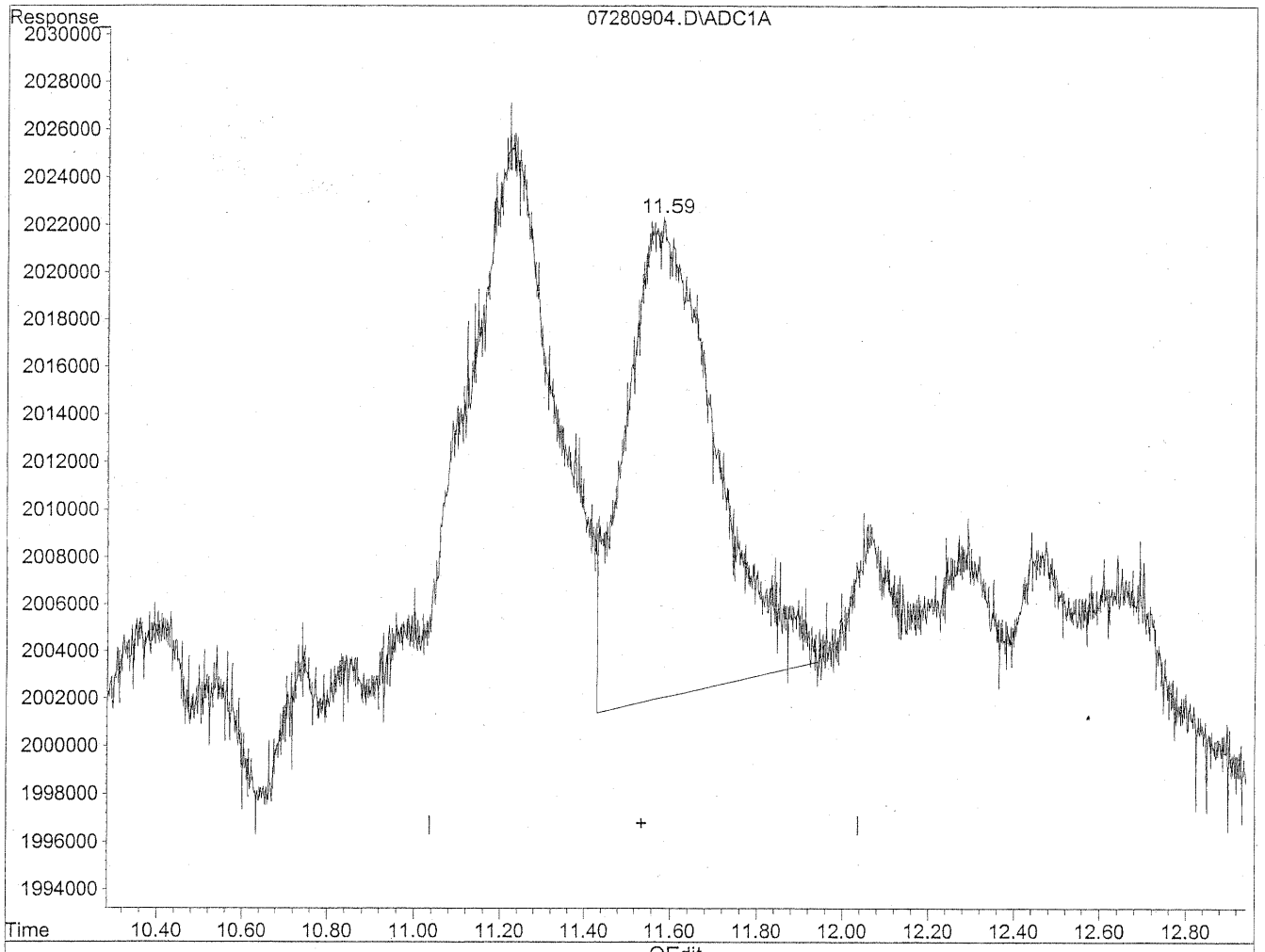
(+) = Expected Retention Time

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



QEdit

(12) 2,5-Dimethylbenzaldehyde

11.58min 55.789ng/ml

response 2897339

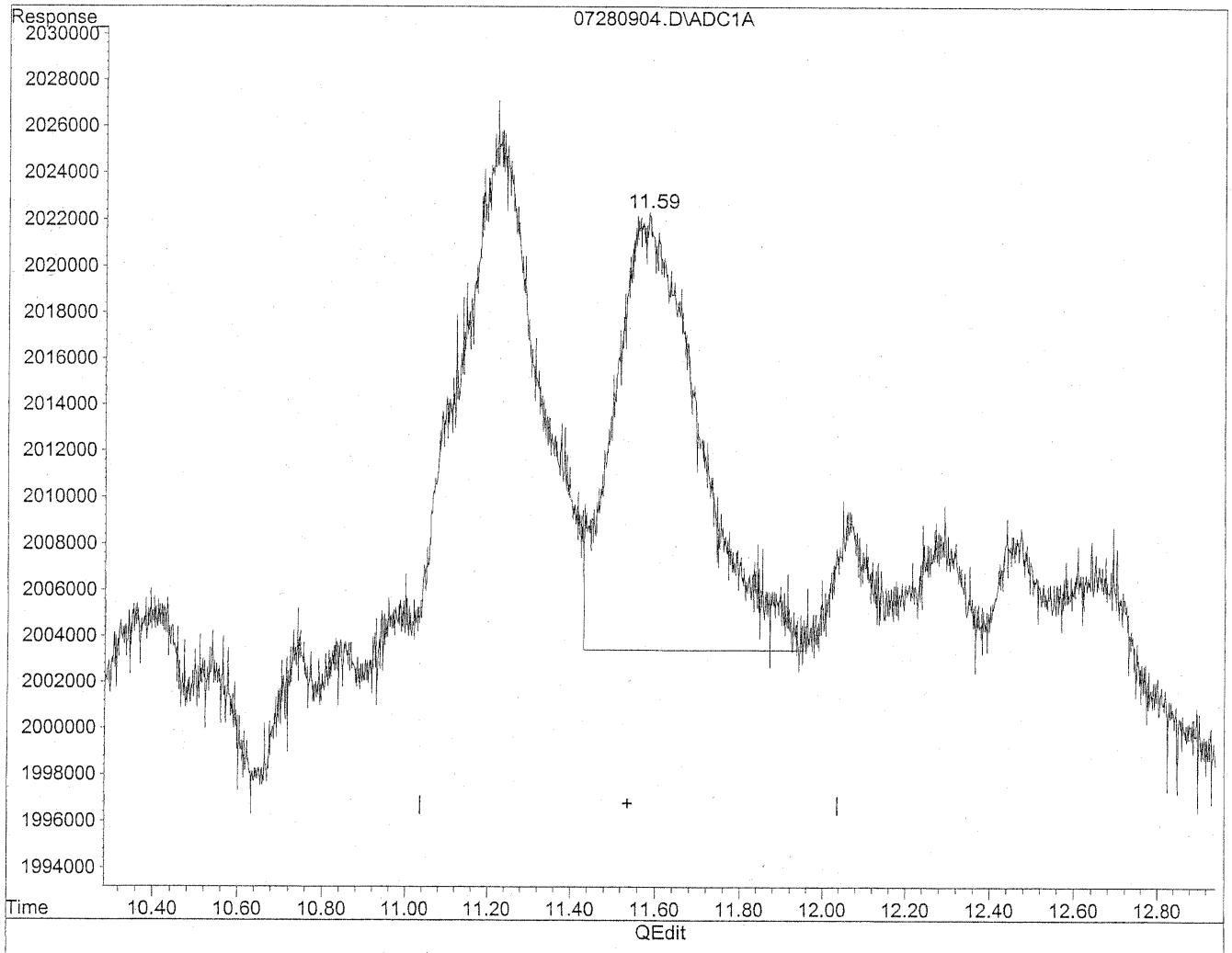
220

(+) = 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

7/28/09

7/29/09

221

(+) = Expected Retention Time

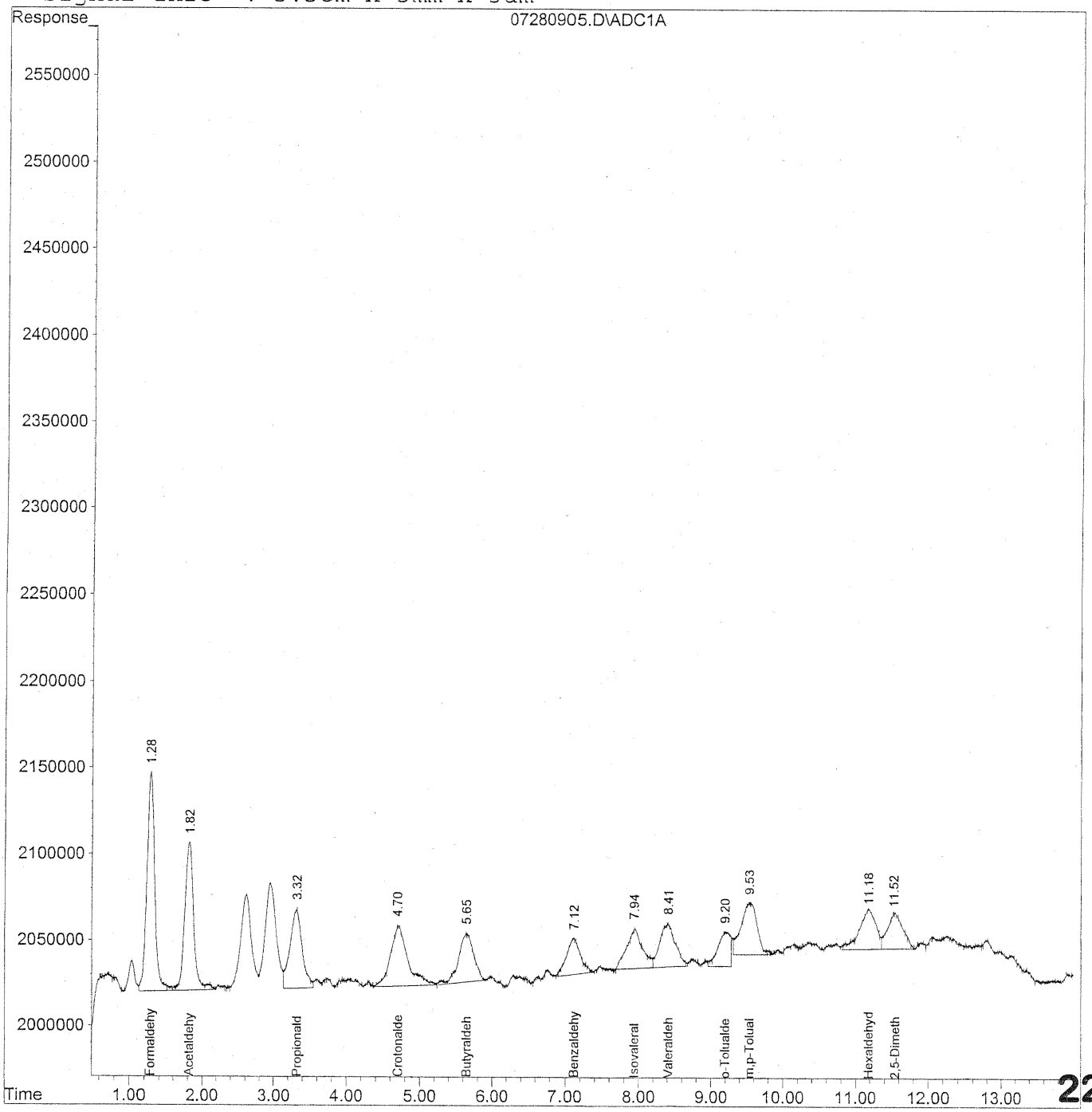
07280904.D TO110709.M Tue Jul 28 14:25:43 2009

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

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



222

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

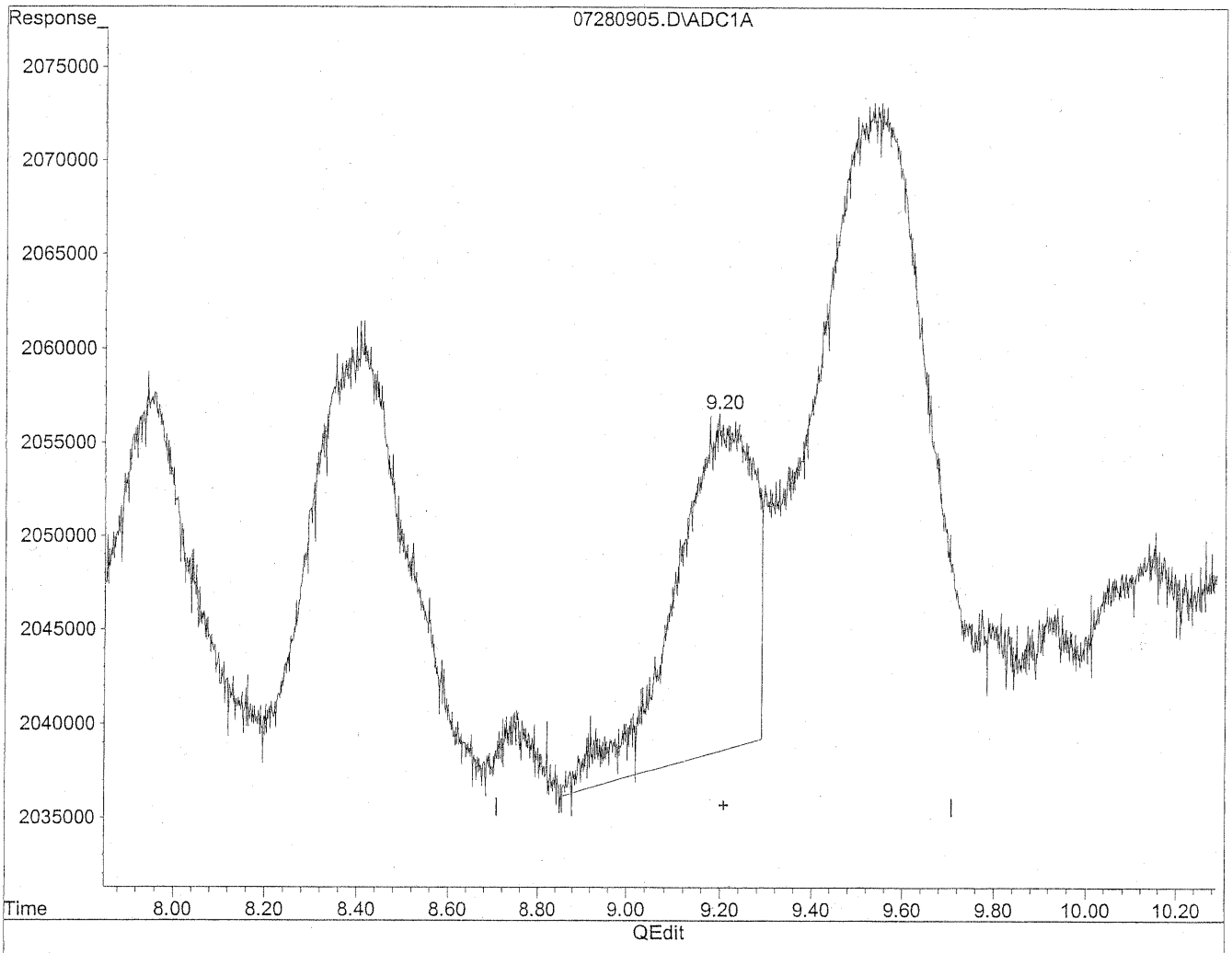
	Compound	R.T.	Response	Conc Units
	Target Compounds			
1)	Formaldehyde	1.28	9305088	52.992 ng/ml
2)	Acetaldehyde	1.81	7389770	54.780 ng/ml
3)	Propionaldehyde	3.31	5442713	52.947 ng/ml
4)	Crotonaldehyde	4.71	5754474	52.051 ng/ml
5)	Butyraldehyde	5.65	4119144	51.185 ng/ml
6)	Benzaldehyde	7.11	2732056	43.316 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

(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

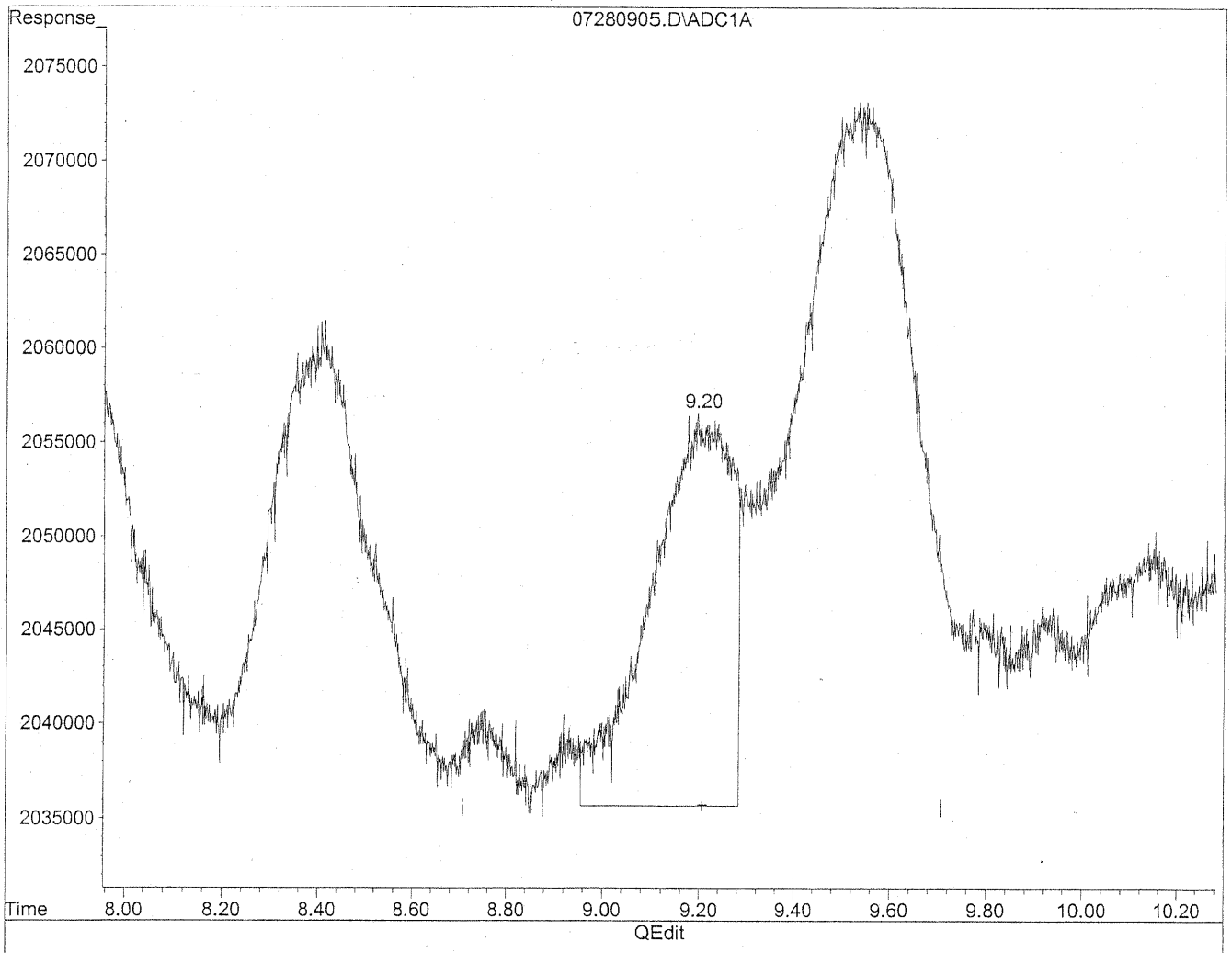
224

(+) = 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 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



(9) o-Tolualdehyde

9.20min 44.856ng/ml m

response 2416389

AL 4/29/09 IC

KE 7/29/09

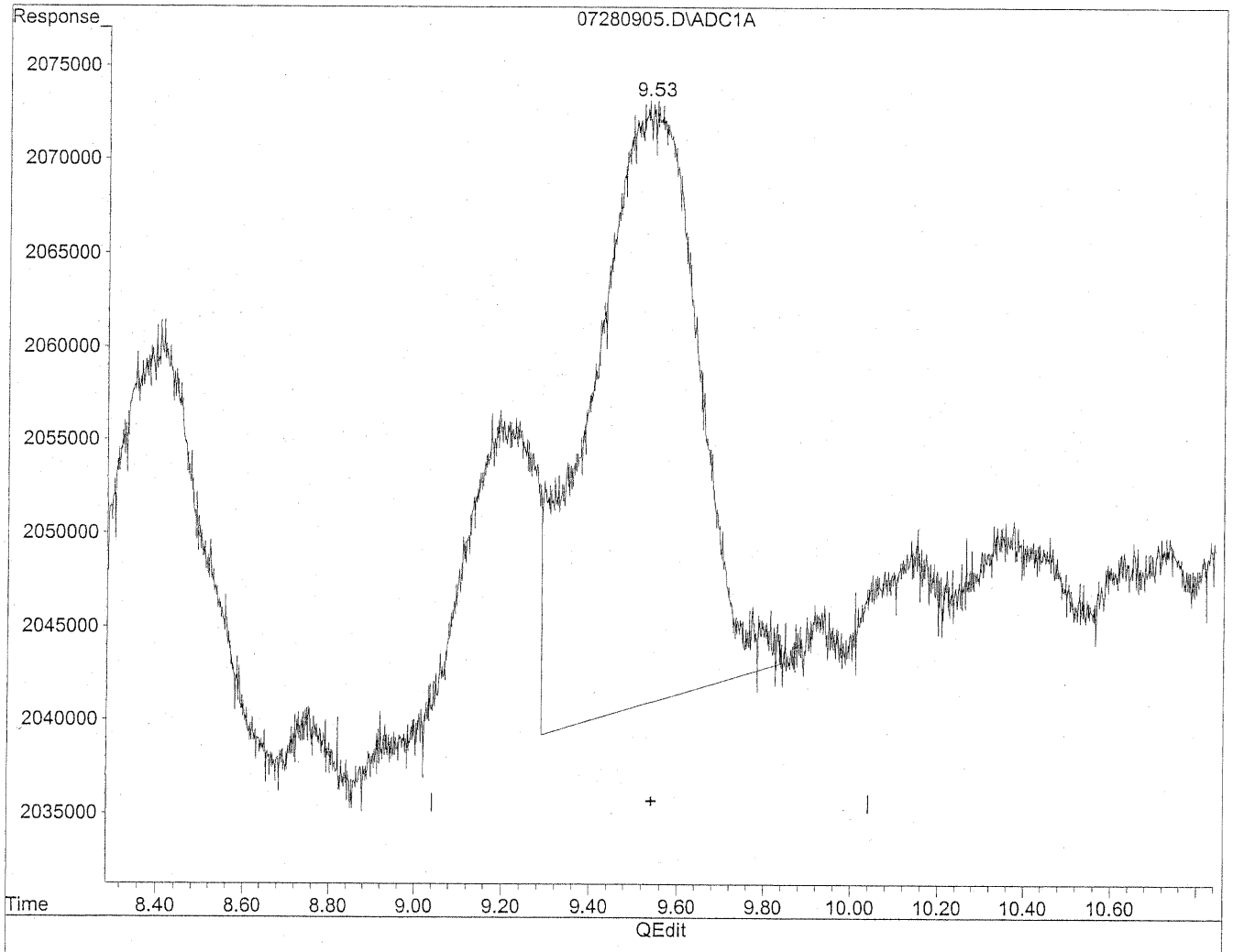
225

(+) = 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

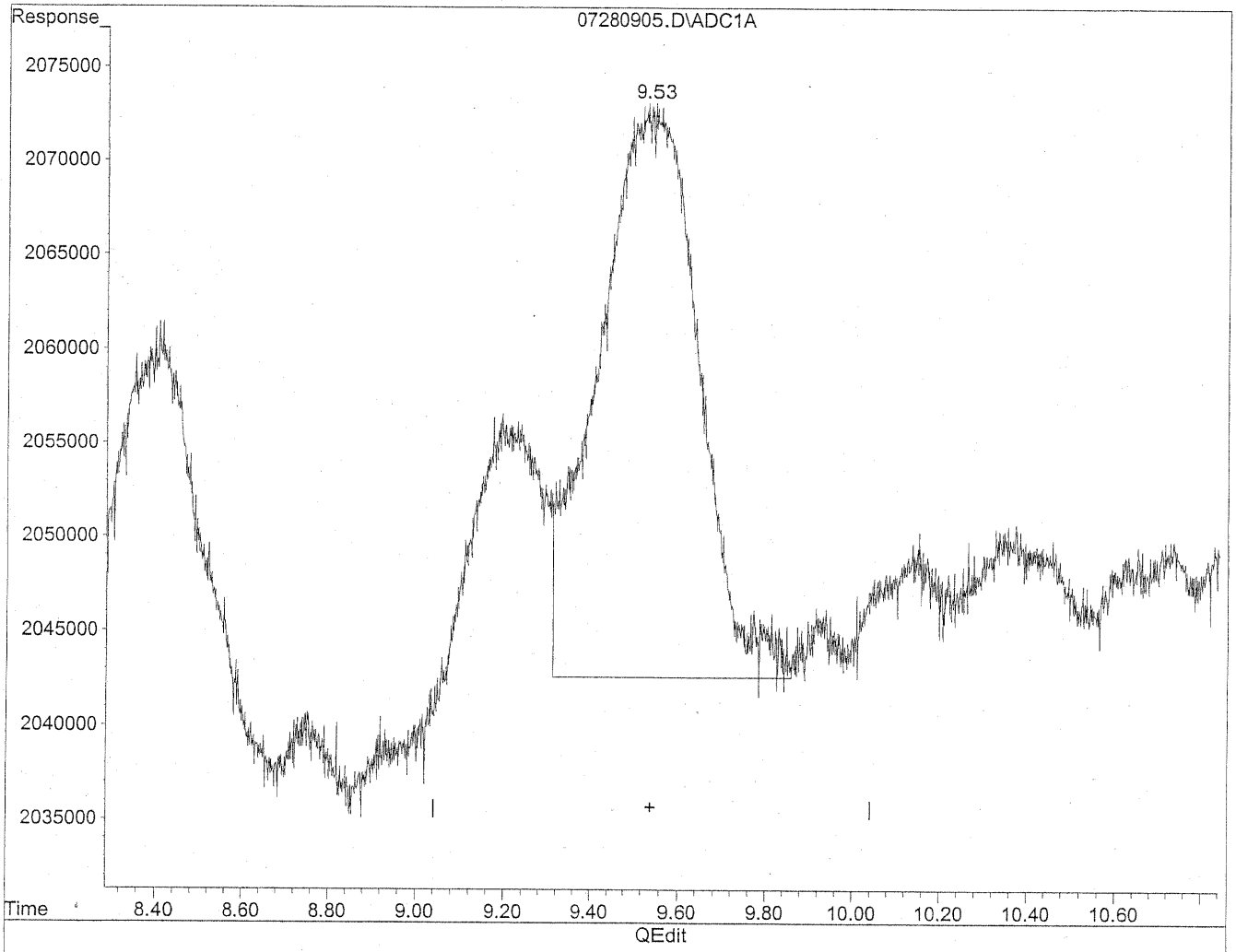
226

(+) = 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 7/28/09 BC

7/29/09

227

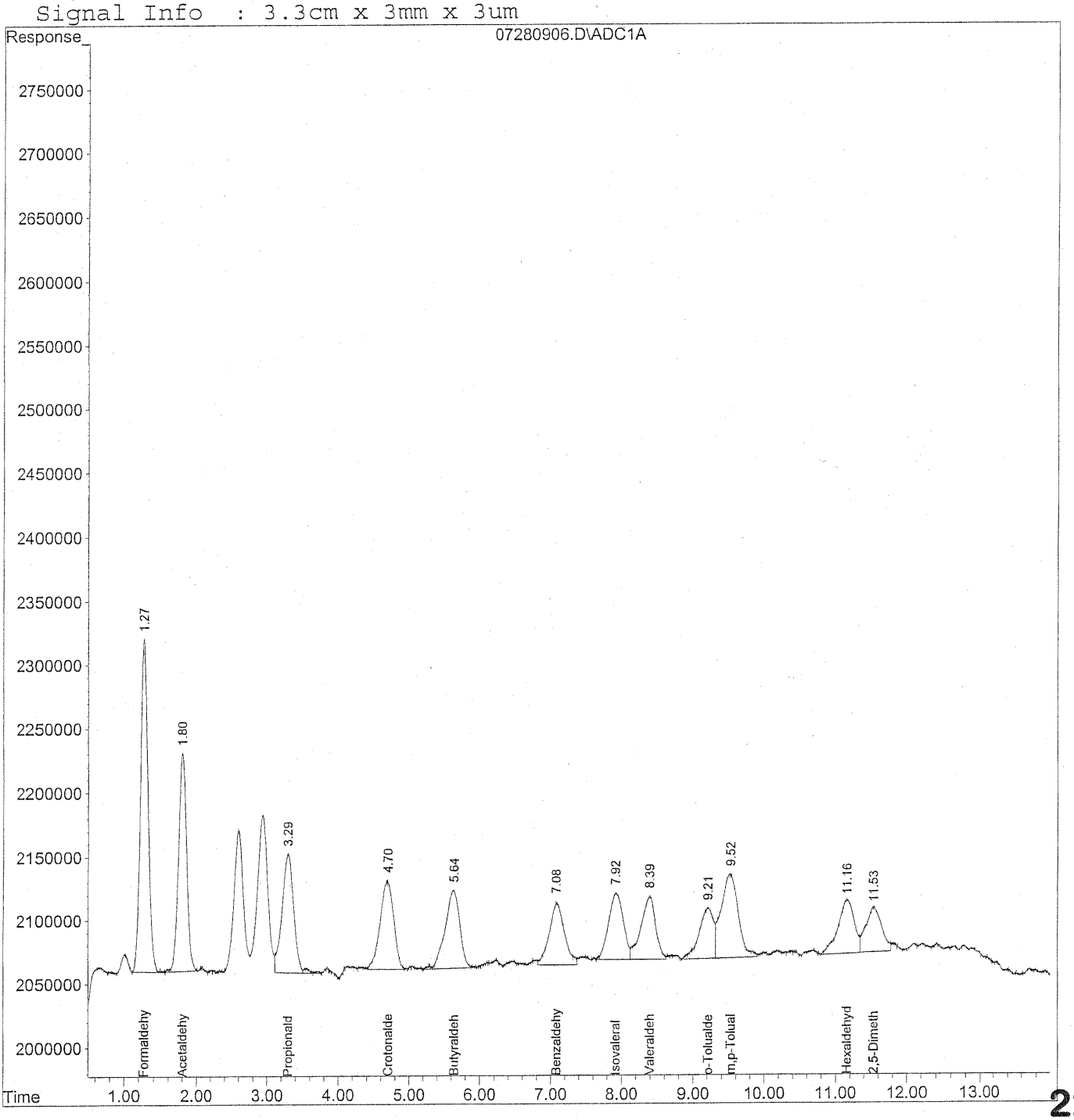
(+) = Expected Retention Time

07280905.D TO110709.M Tue Jul 28 14:29:37 2009

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: Sep 10 9:16 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 : 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



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: Sep 10 9:16 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 : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

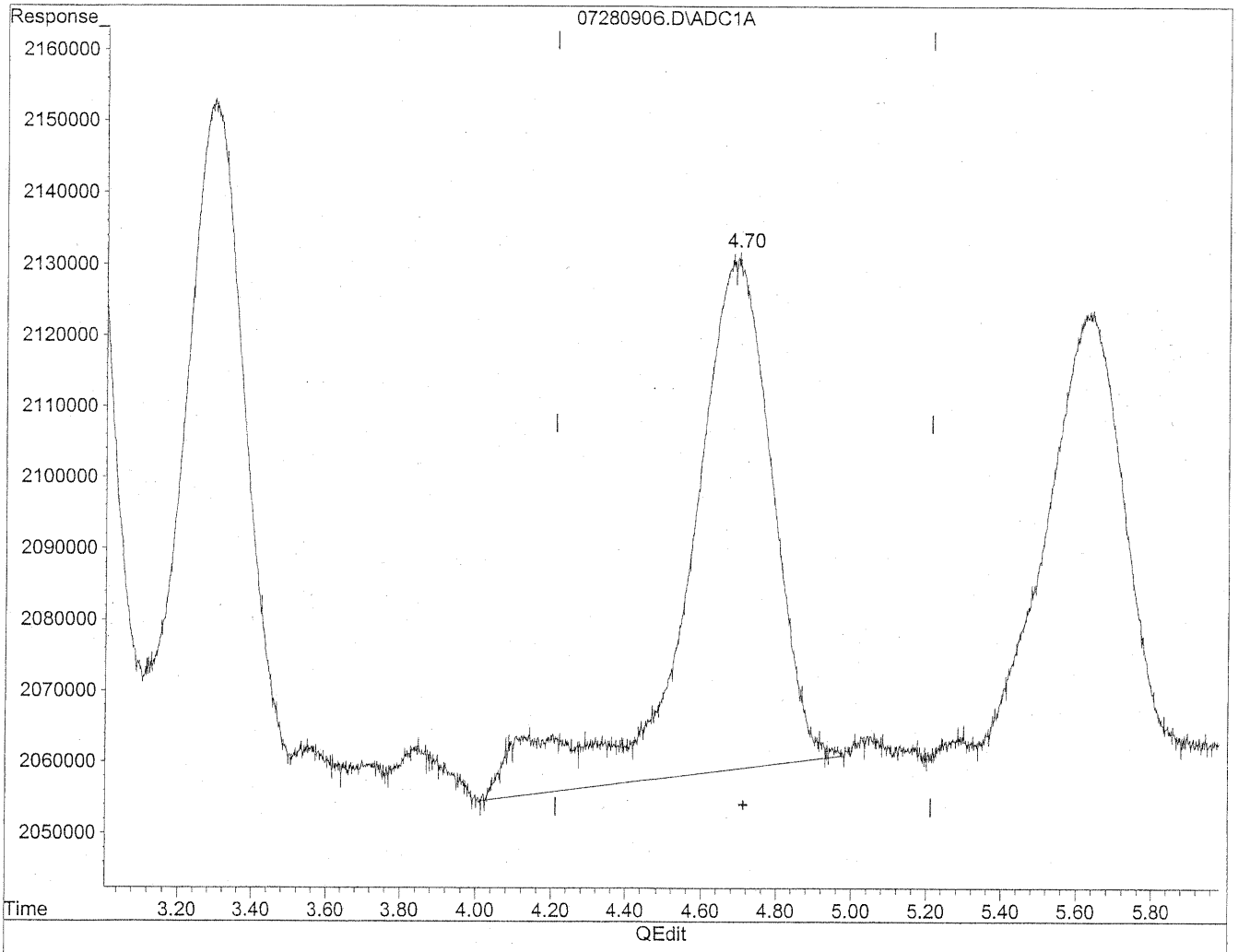
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	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	5399082	103.961	ng/mlm

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

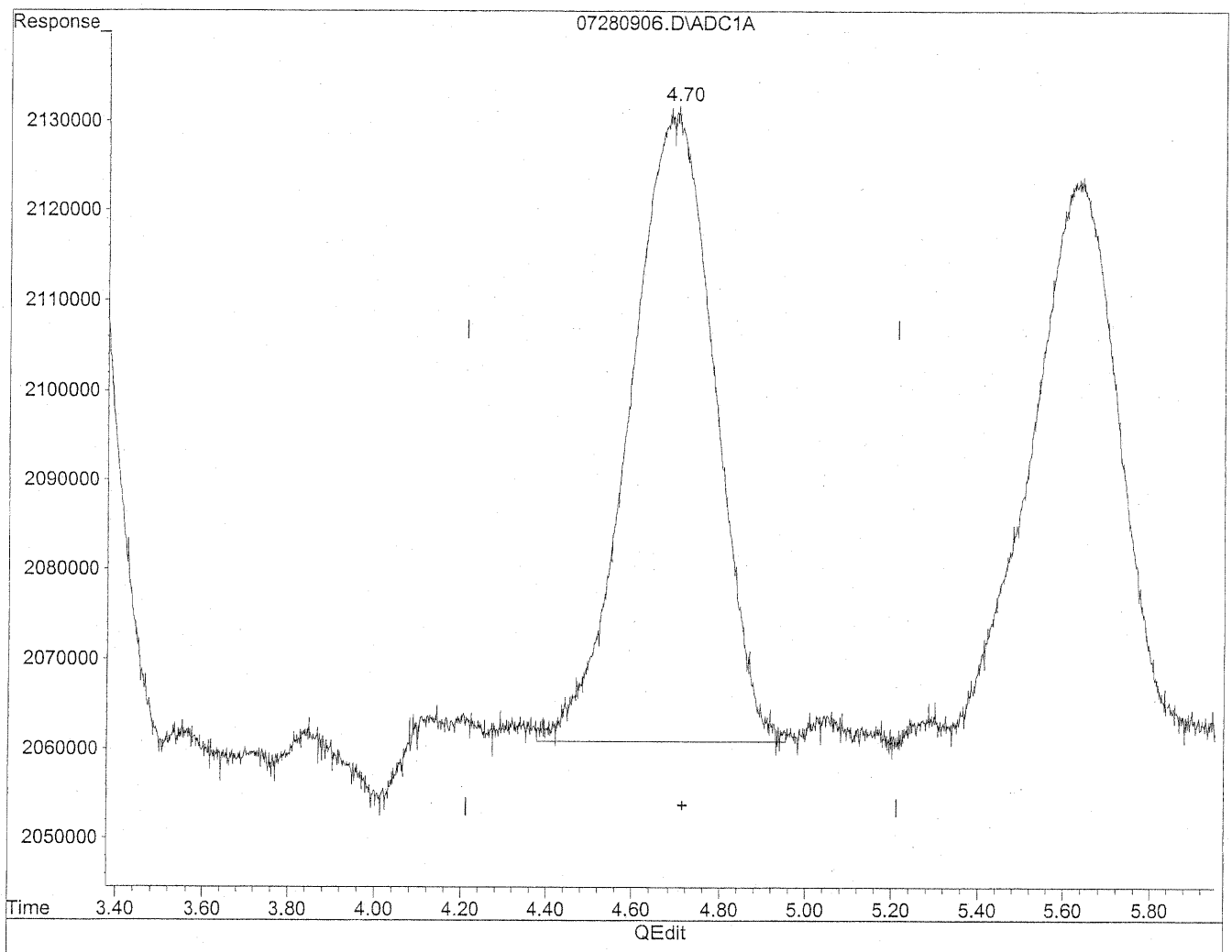
response 11317409

(+) = 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

JLC 07/28/09 LC

KW 7/29/09

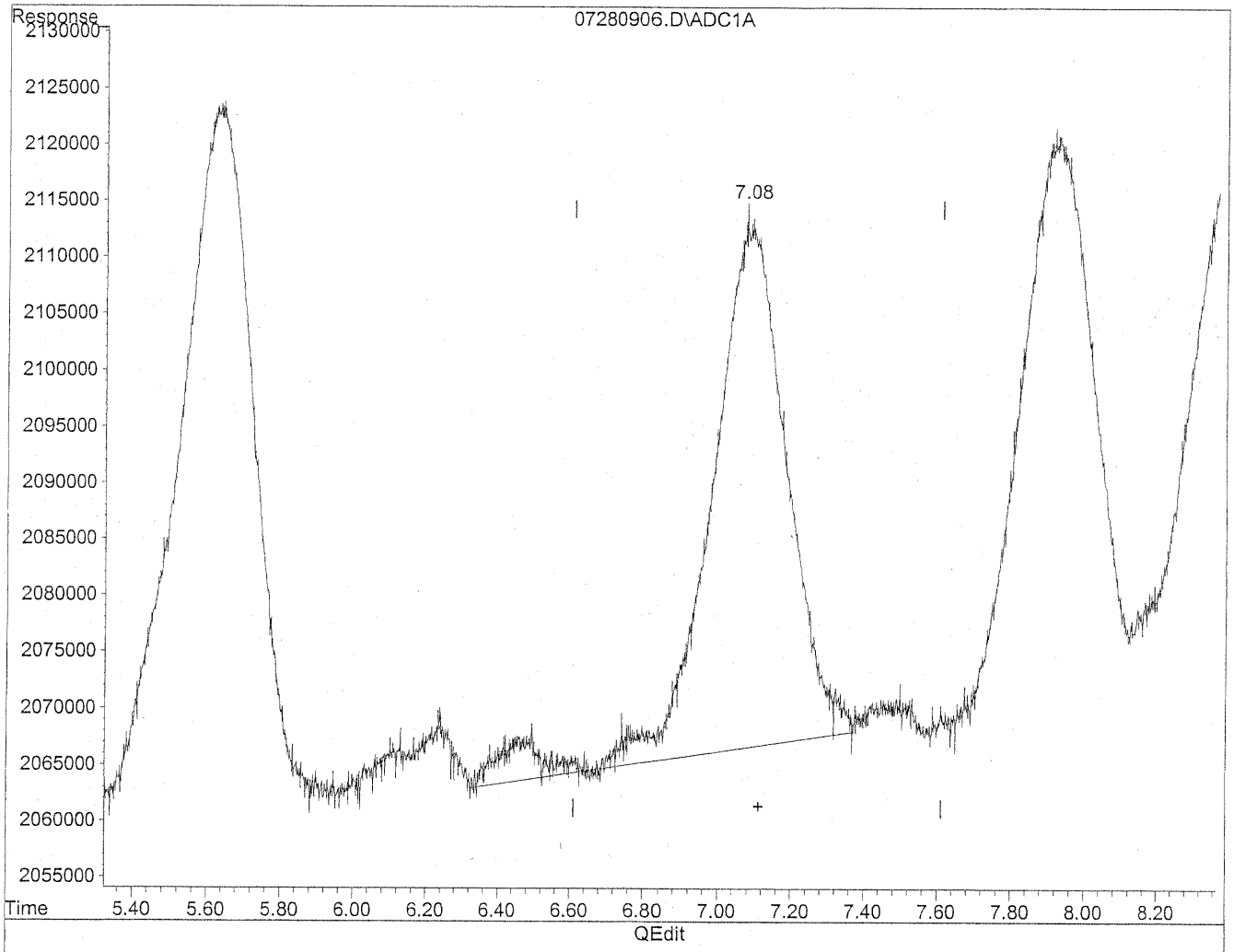
231

(+) = 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

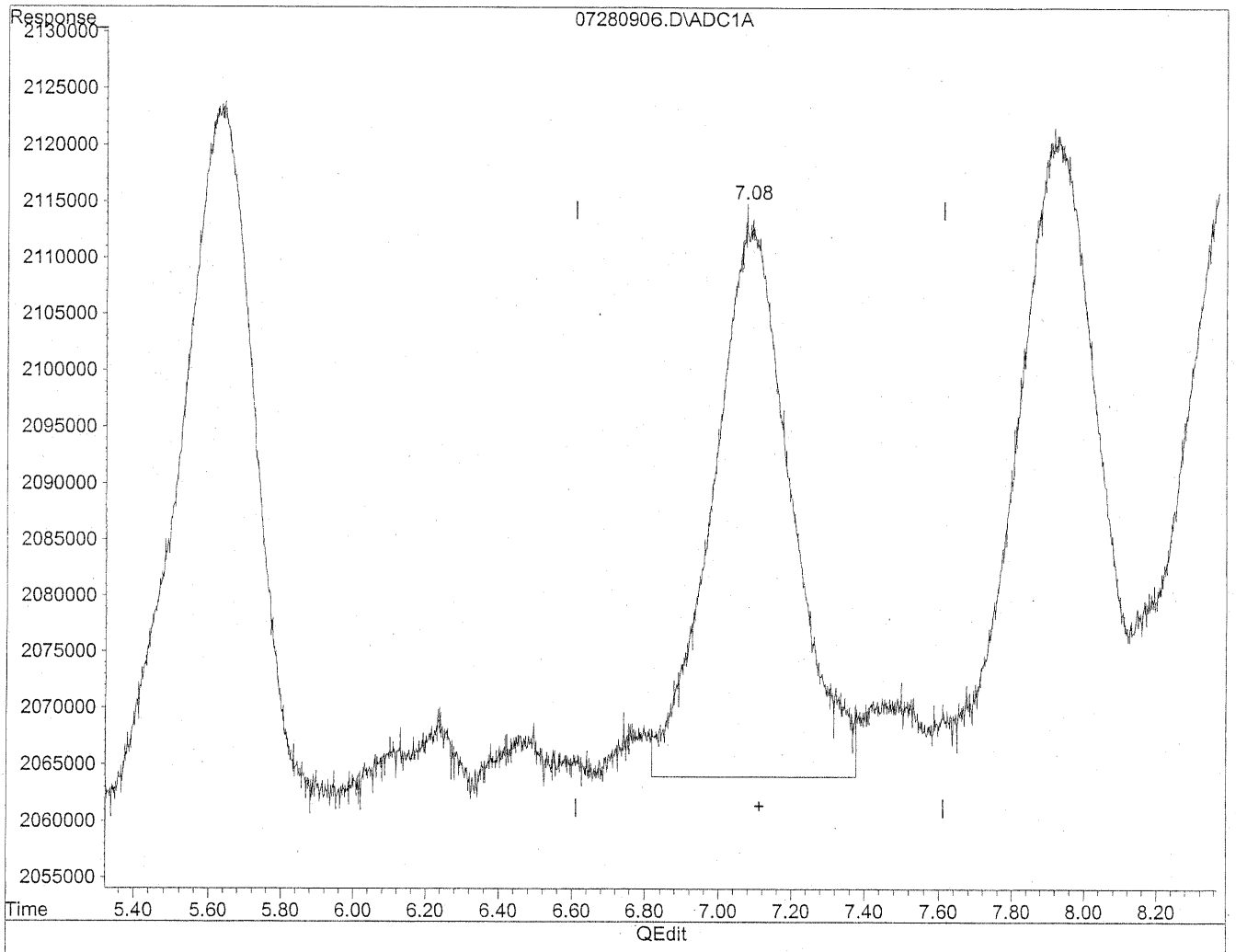
response 6819663

(+) = 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

7/28/09 JC

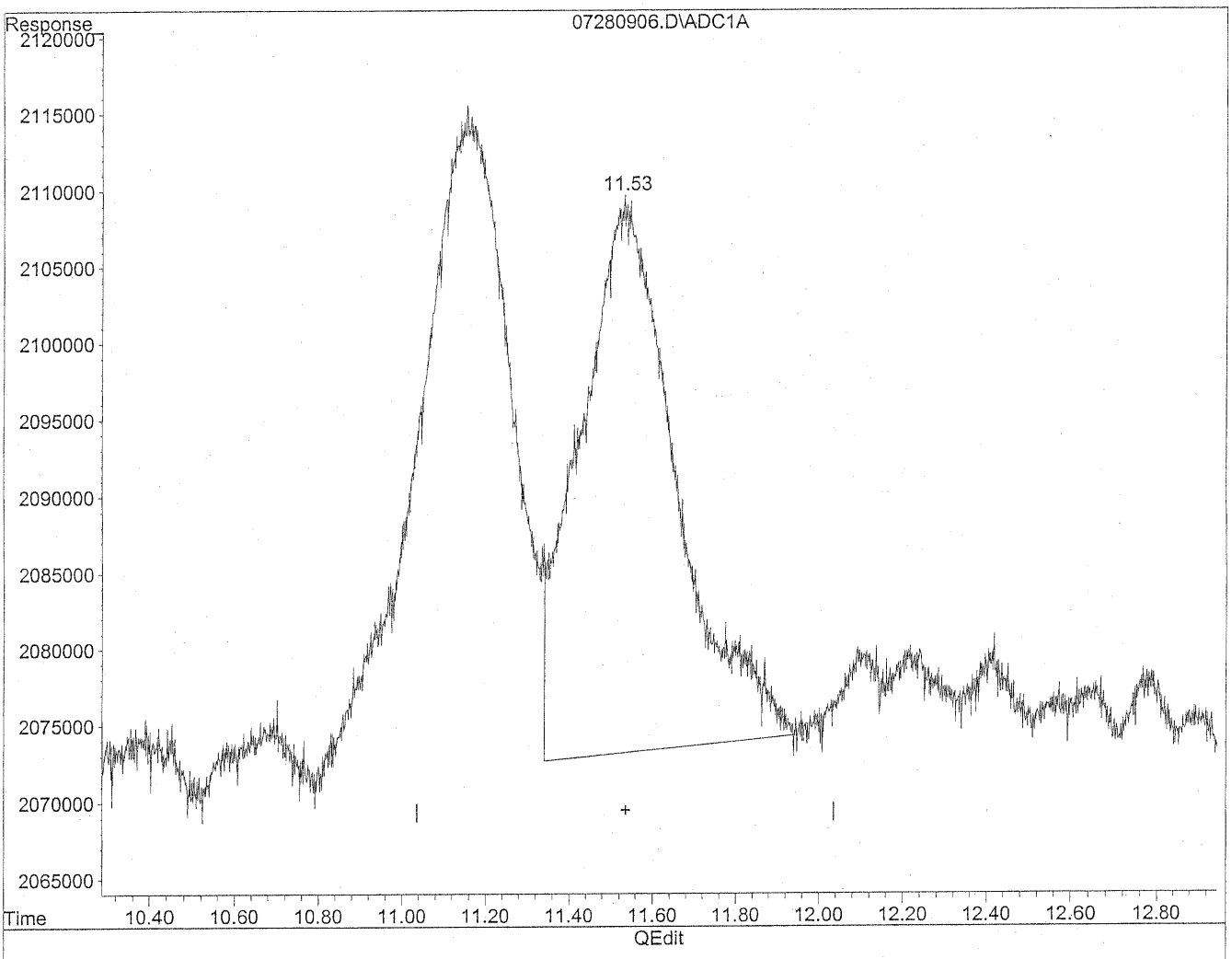
KR 7/29/09

233

(+) = Expected Retention Time

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 : Mon Mar 21 12:19:47 2005
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

11.53min 111.652ng/ml

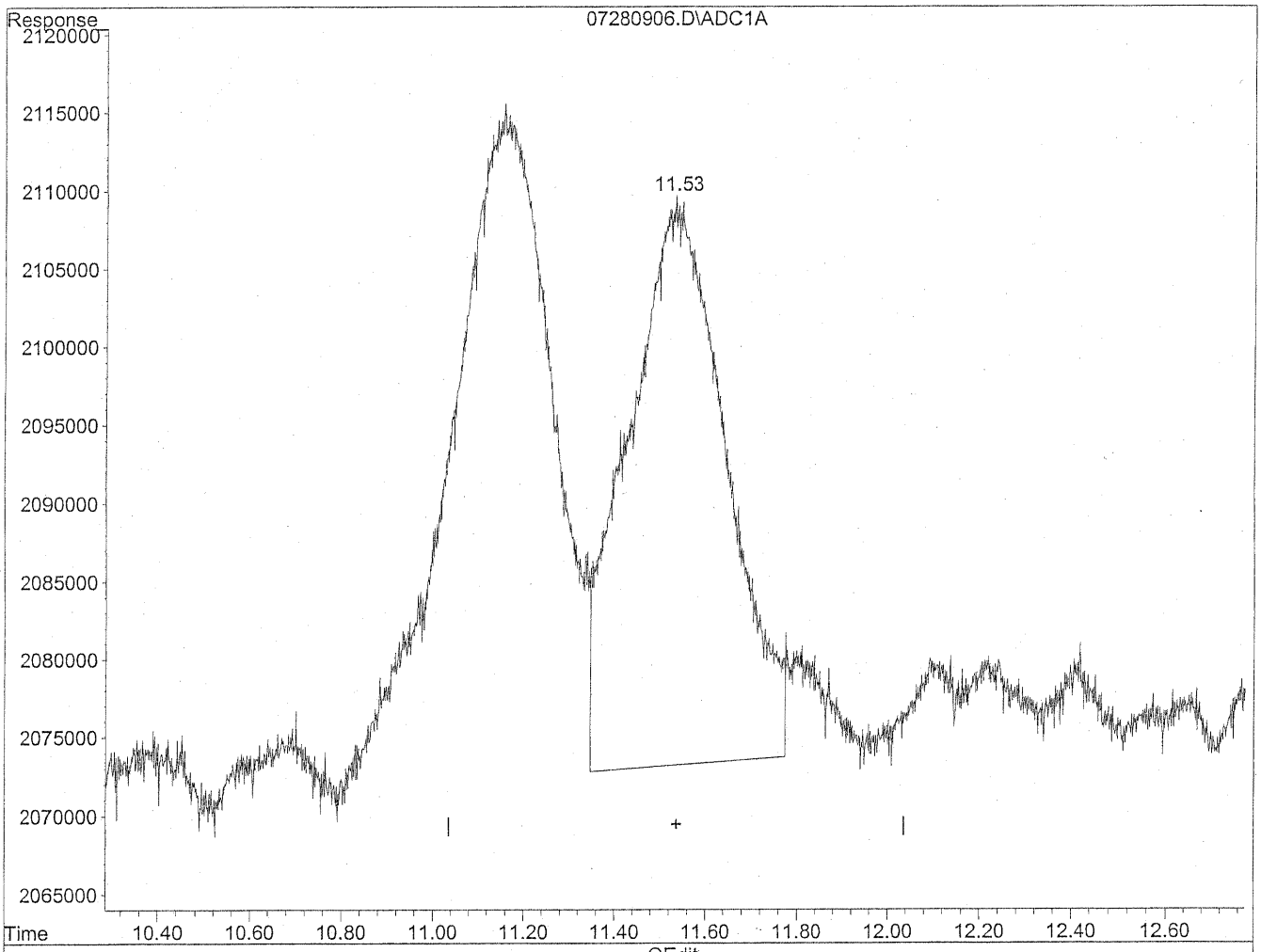
response 5798505

(+) = 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: Sep 10 9:16 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO11709B.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

11.53min 103.961ng/ml m

response 5399082

HL & 09/10/09 BC

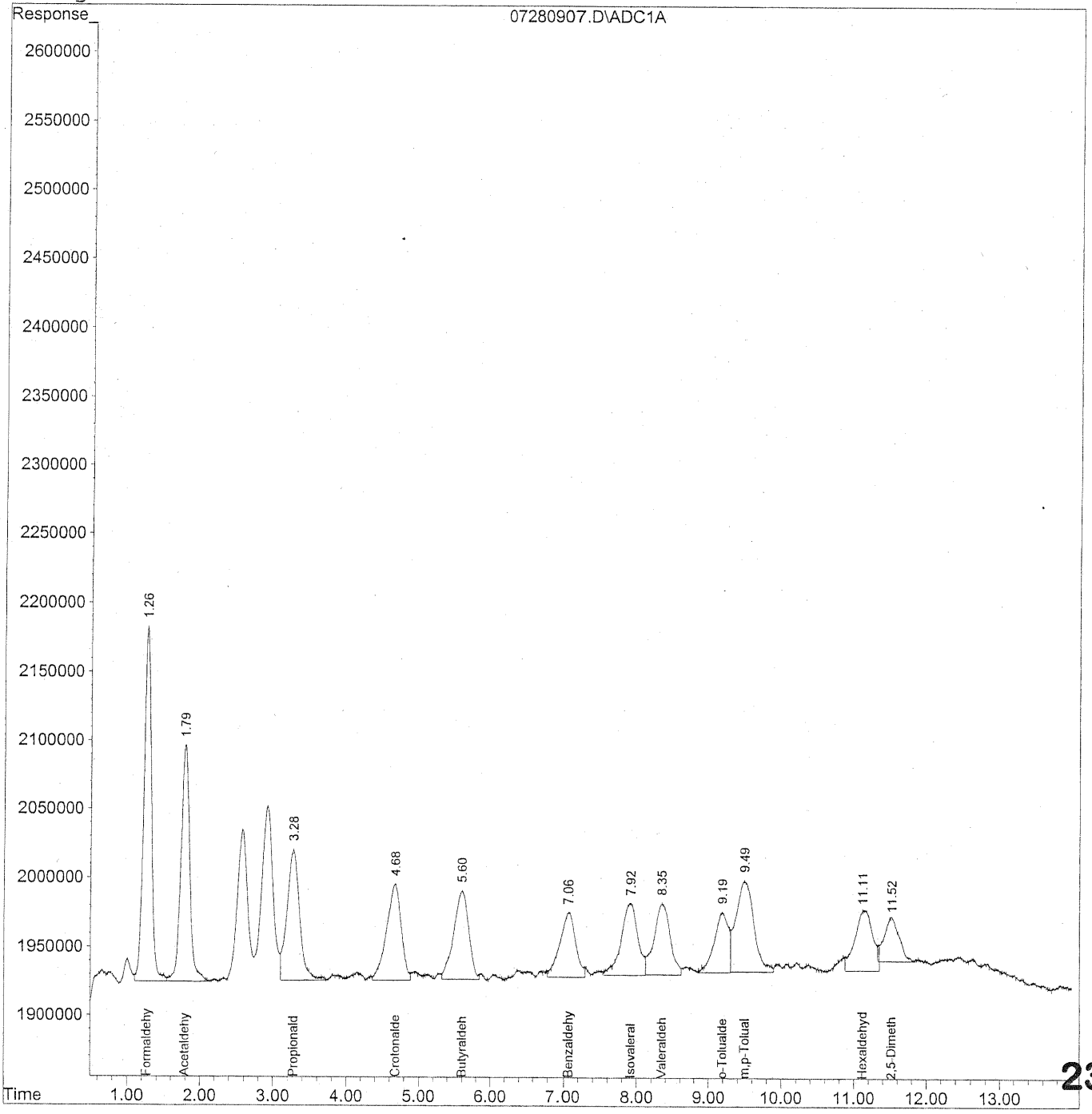
(+) = Expected Retention Time

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 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



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

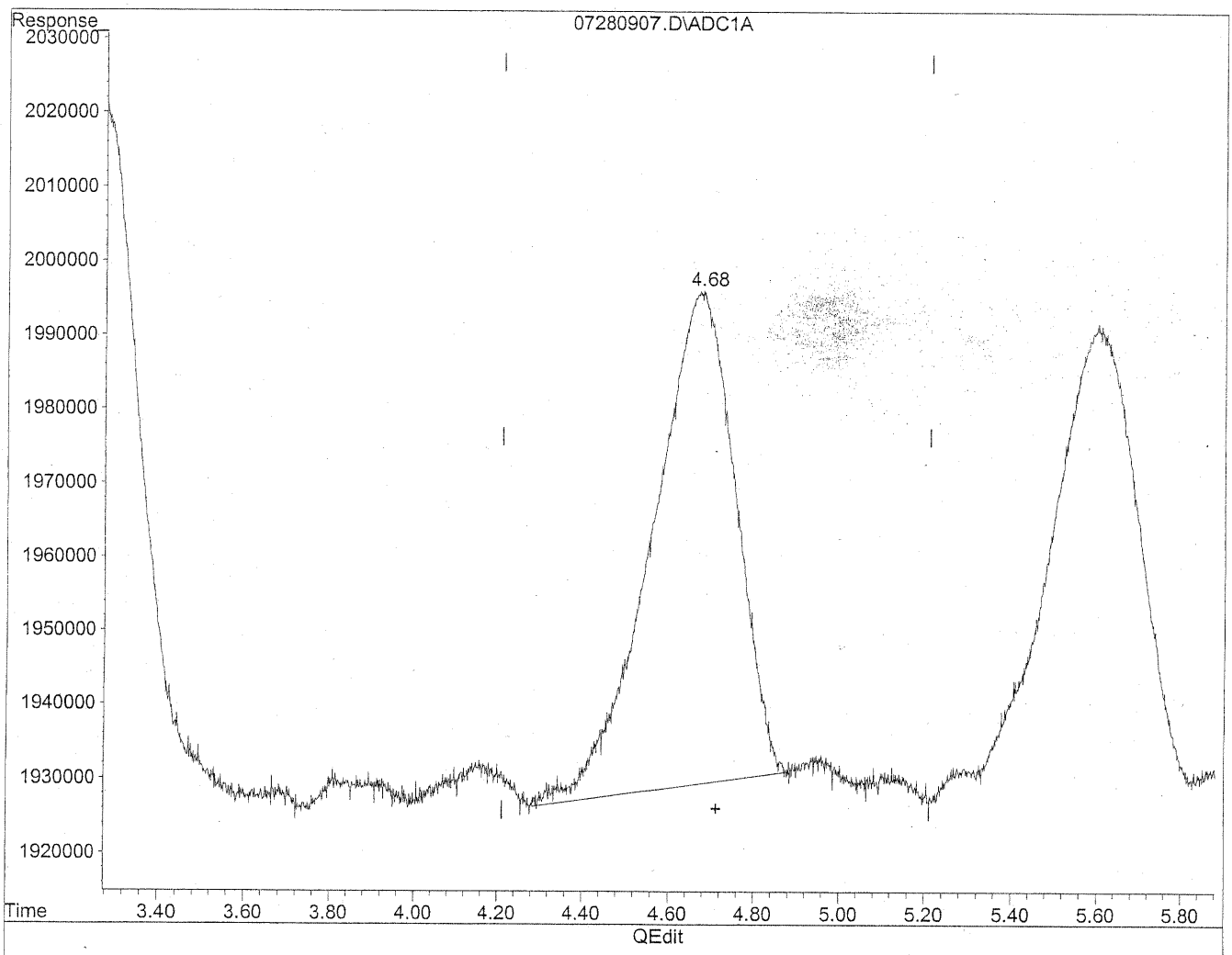
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.26	18449443	105.069	ng/ml
2)	Acetaldehyde	1.79	14434553	107.002	ng/ml
3)	Propionaldehyde	3.28	11389784	110.800	ng/ml
4)	Crotonaldehyde	4.68	9814490	88.774	ng/mlm
5)	Butyraldehyde	5.60	9432197	117.206	ng/mlm
6)	Benzaldehyde	7.06	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.

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

238

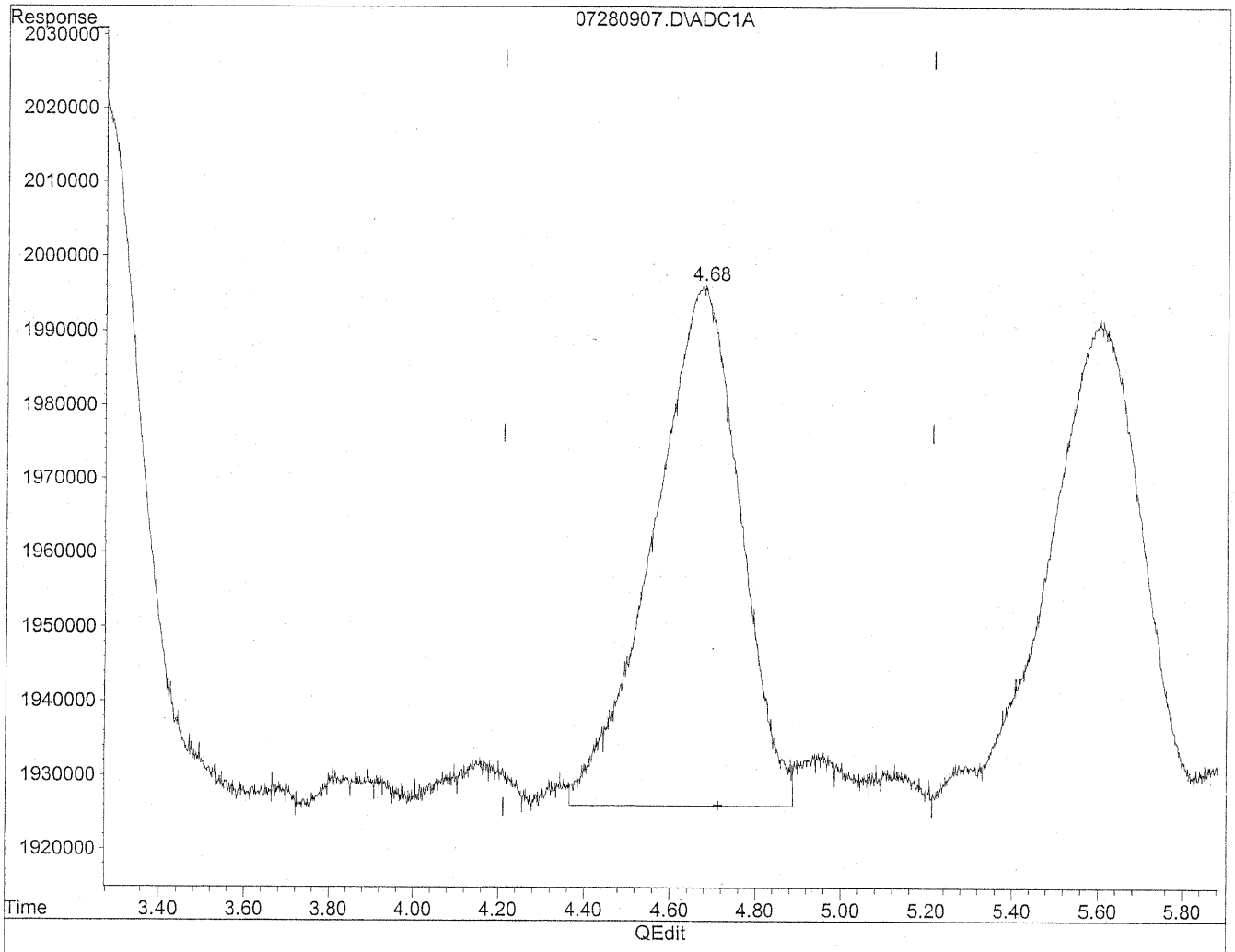
(+) = Expected Retention Time

07280907.D TO110709.M Tue Jul 28 14:31:01 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

4.68min 88.774ng/ml m

response 9814490

AL 7/28/09 1C

KR 7/29/09

239

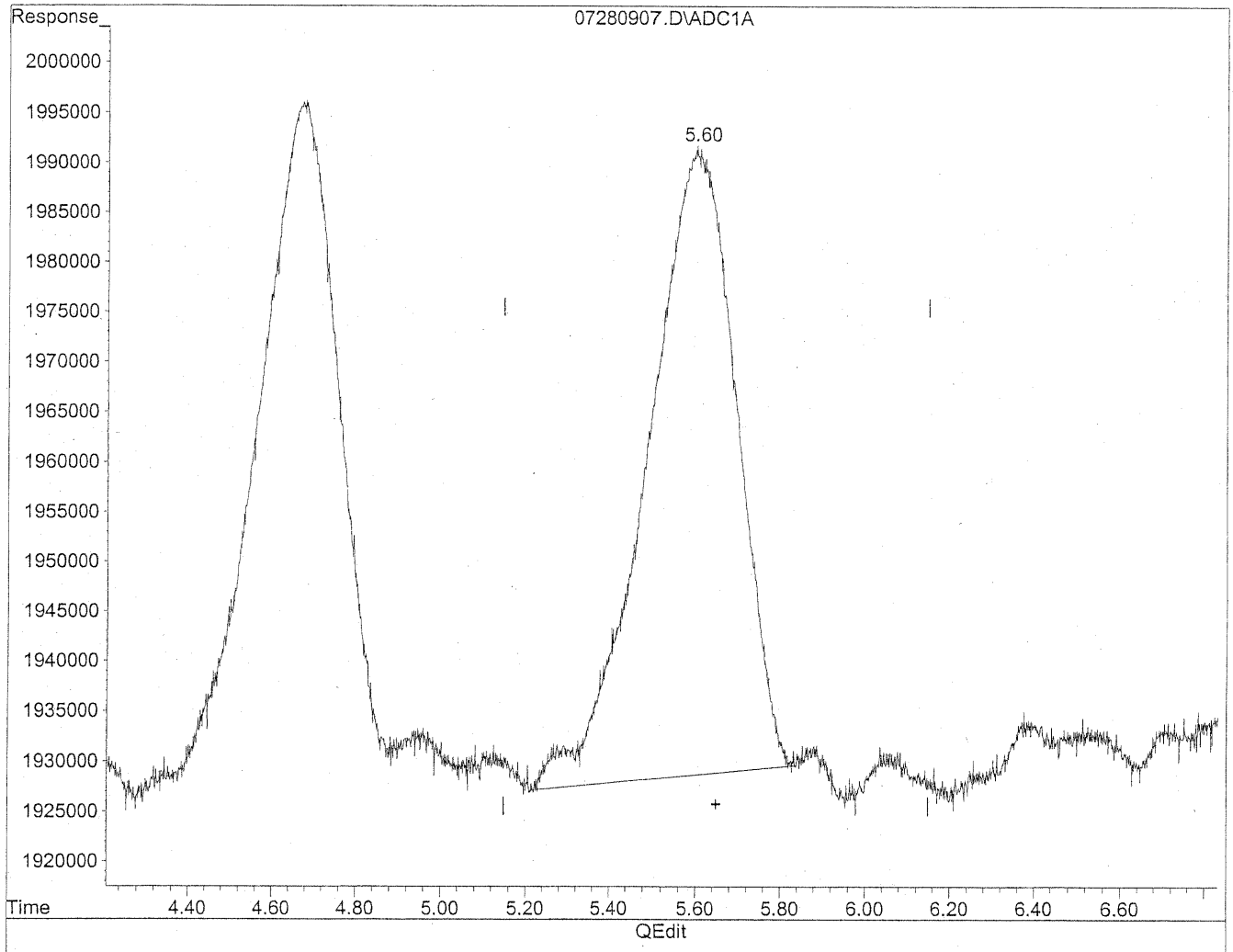
(+) = Expected Retention Time

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

240

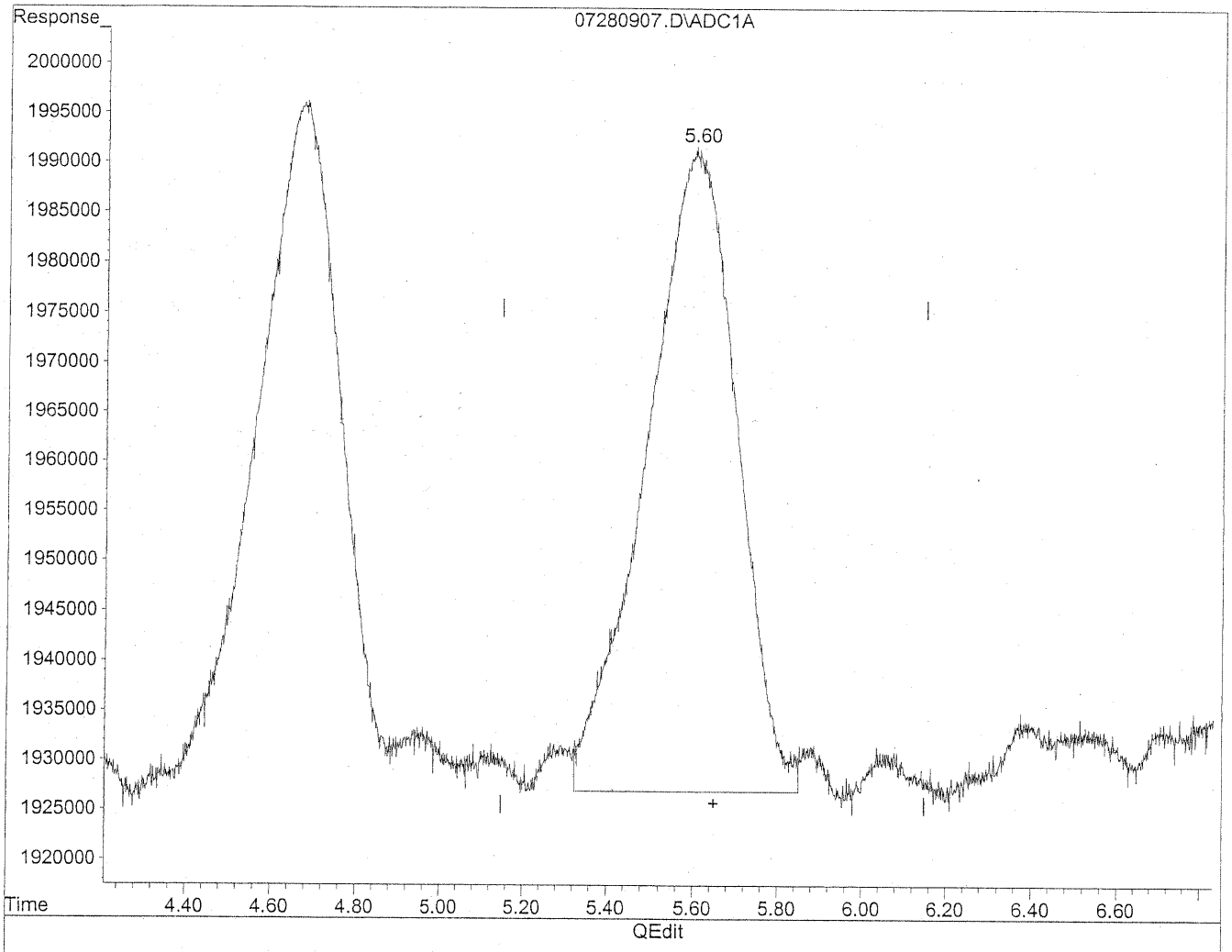
(+) = Expected Retention Time

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 7/28/09 LC

KE 7/29/09

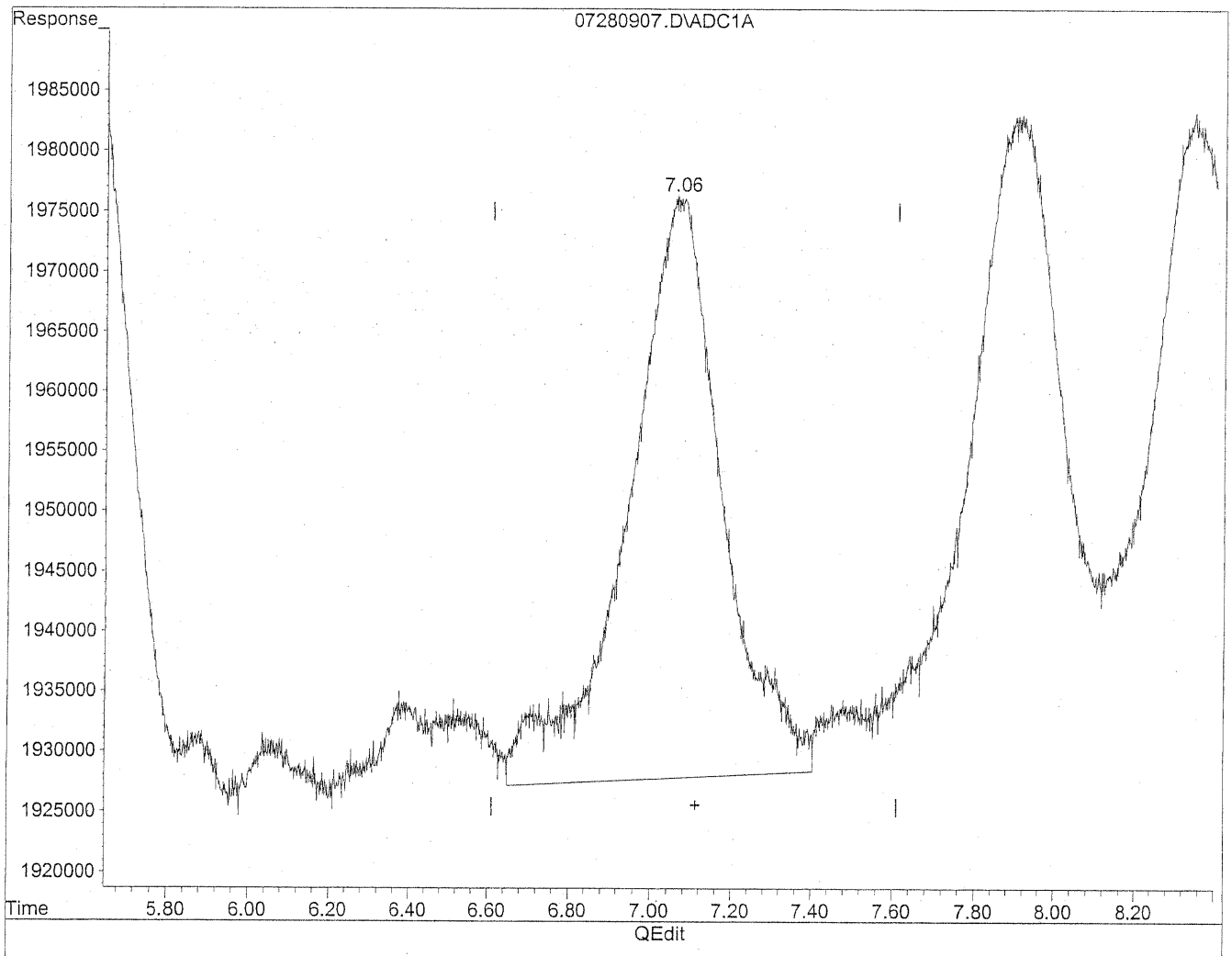
241

(+) = 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

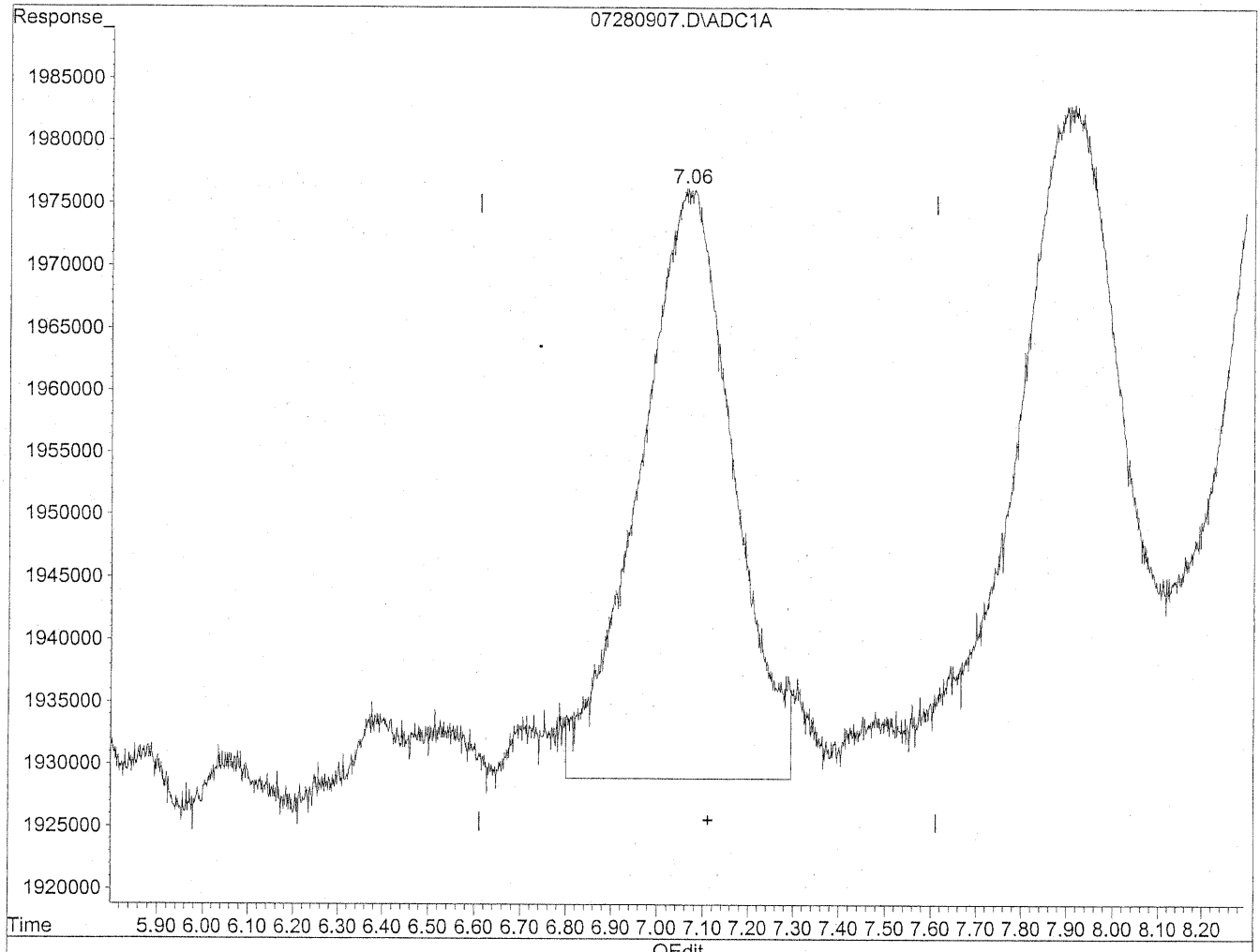
response 7772036

(+) = Expected Retention Time

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

KR 7/29/09

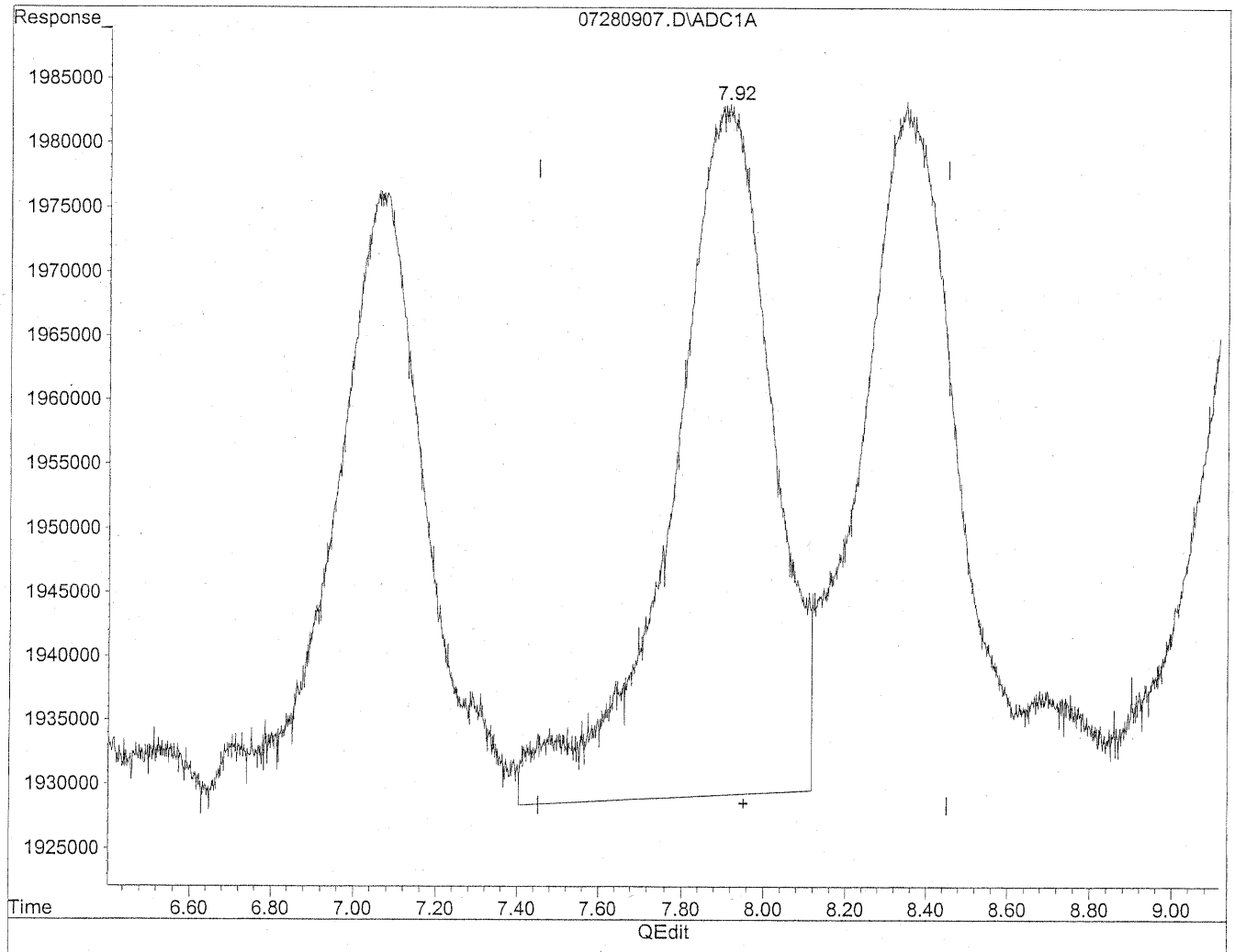
243

(+) = 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

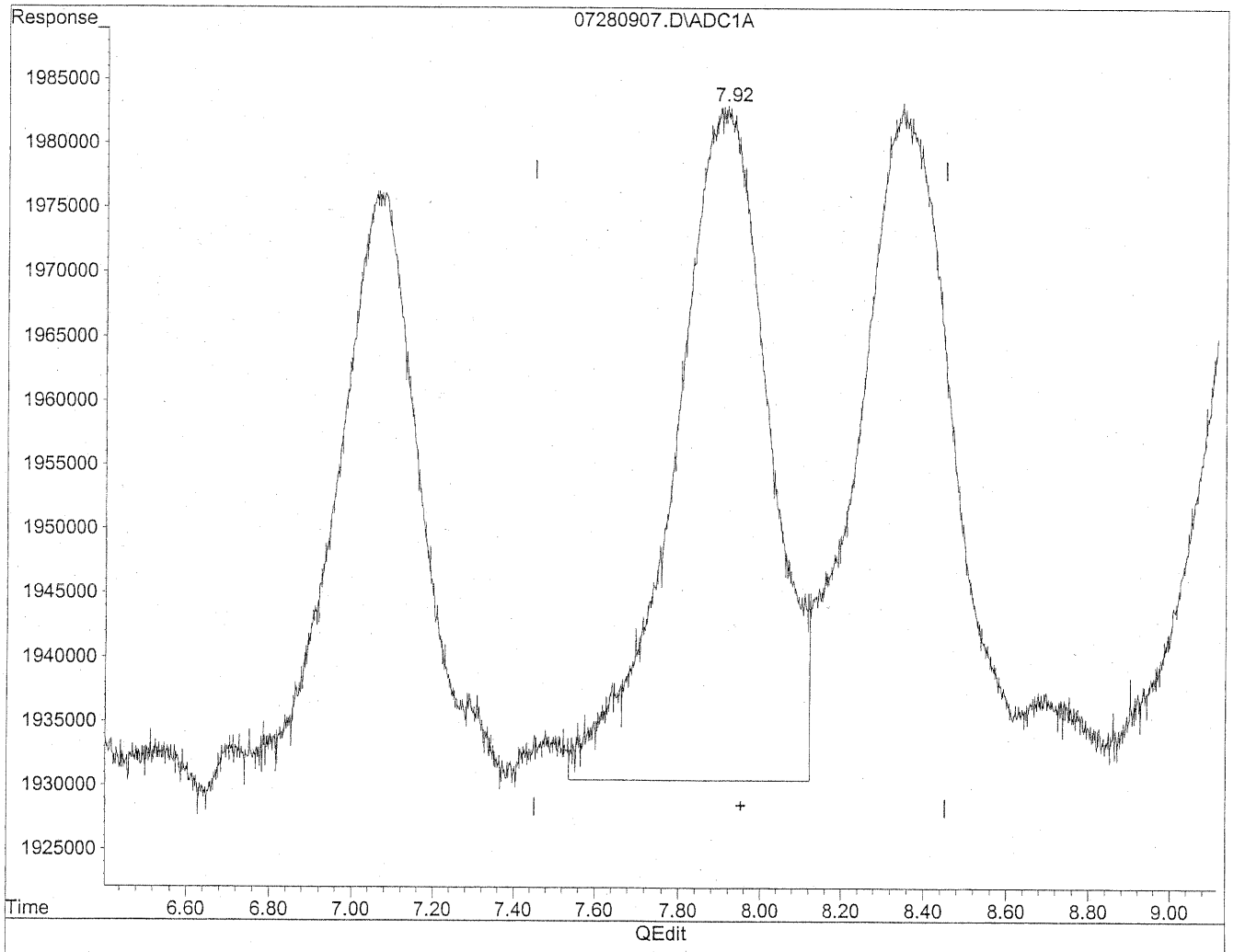
response 9140643

(+) = 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

HL 7/28/09 IC

KR 7/29/09

245

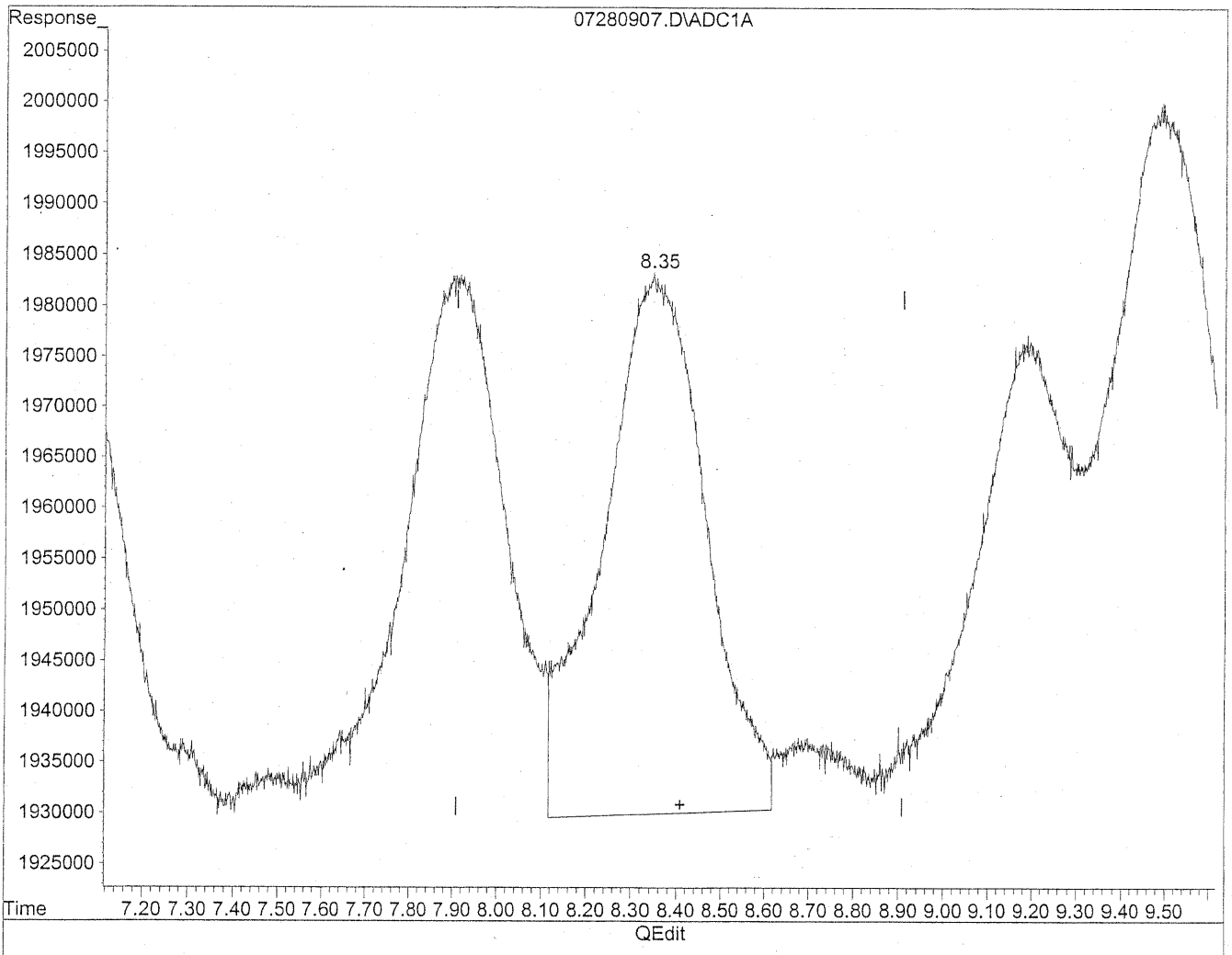
(+) = Expected Retention Time

07280907.D TO110709.M Tue Jul 28 14:31:57 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



(8) Valeraldehyde

8.35min 101.373ng/ml

response 8423554

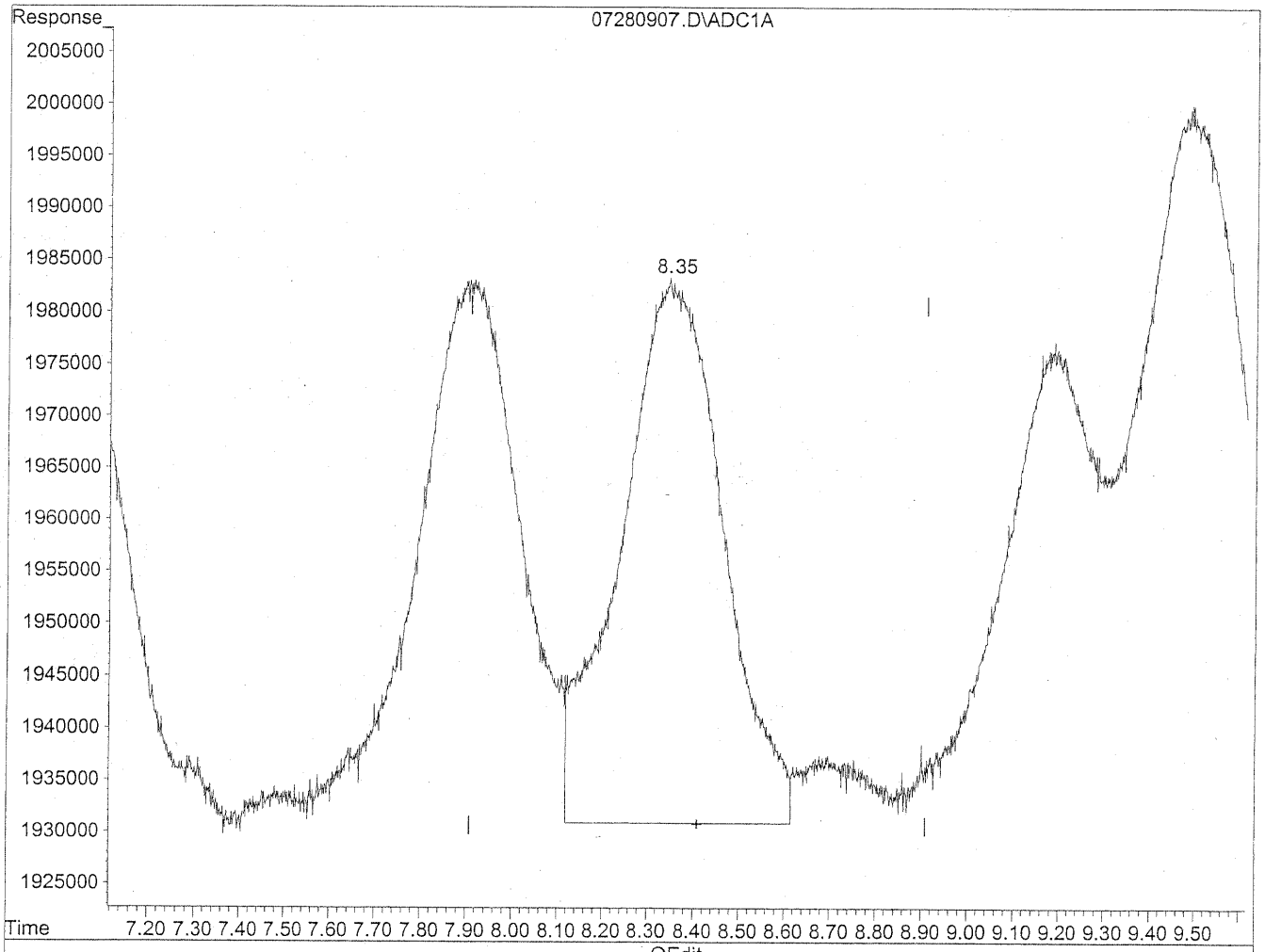
246

(+) = 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 97.688ng/ml m

response 8117341

HC 7/28/09 RC

KR 7/29/09

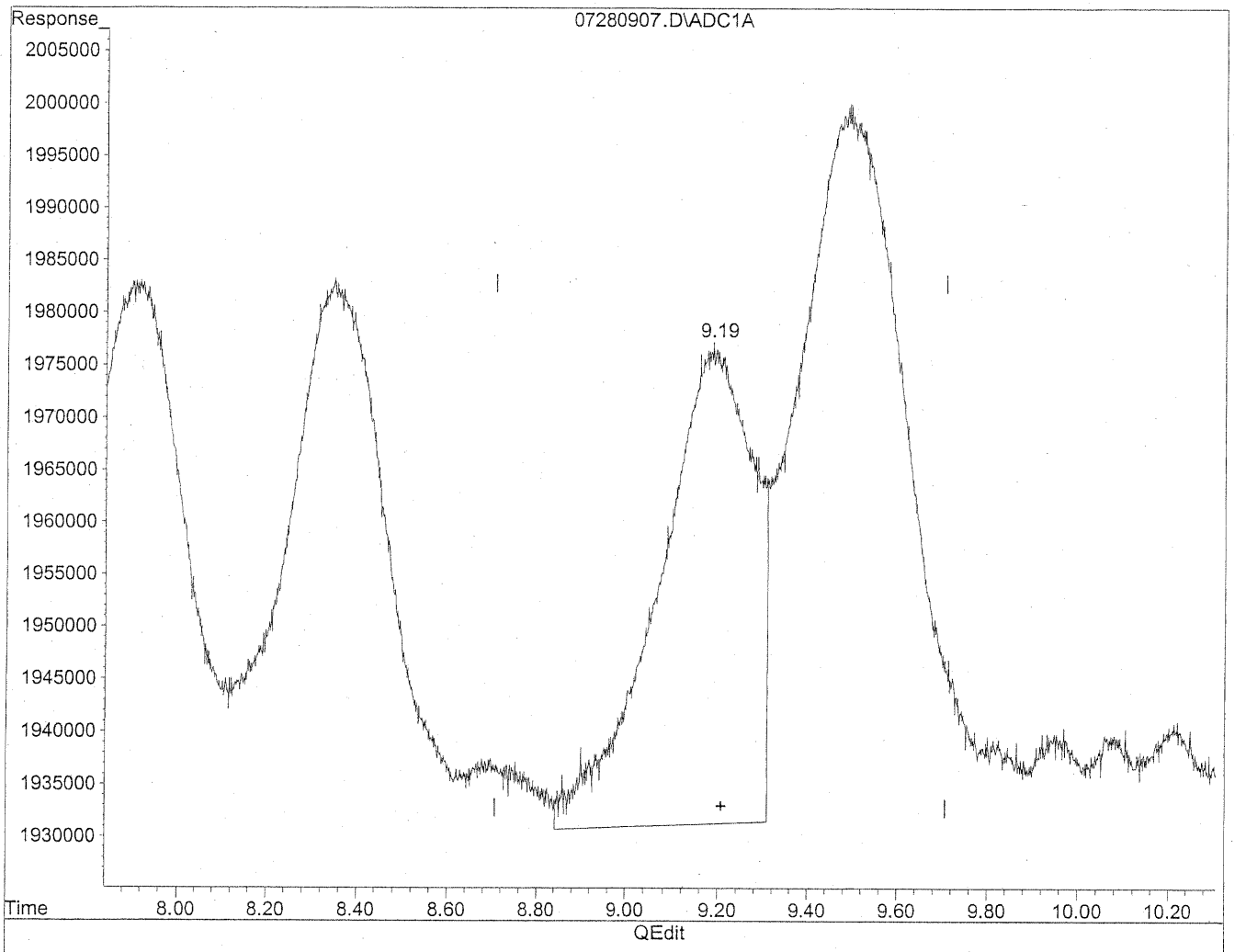
247

(+) = 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



(9) o-Tolualdehyde

9.19min 121.312ng/ml

response 6535124

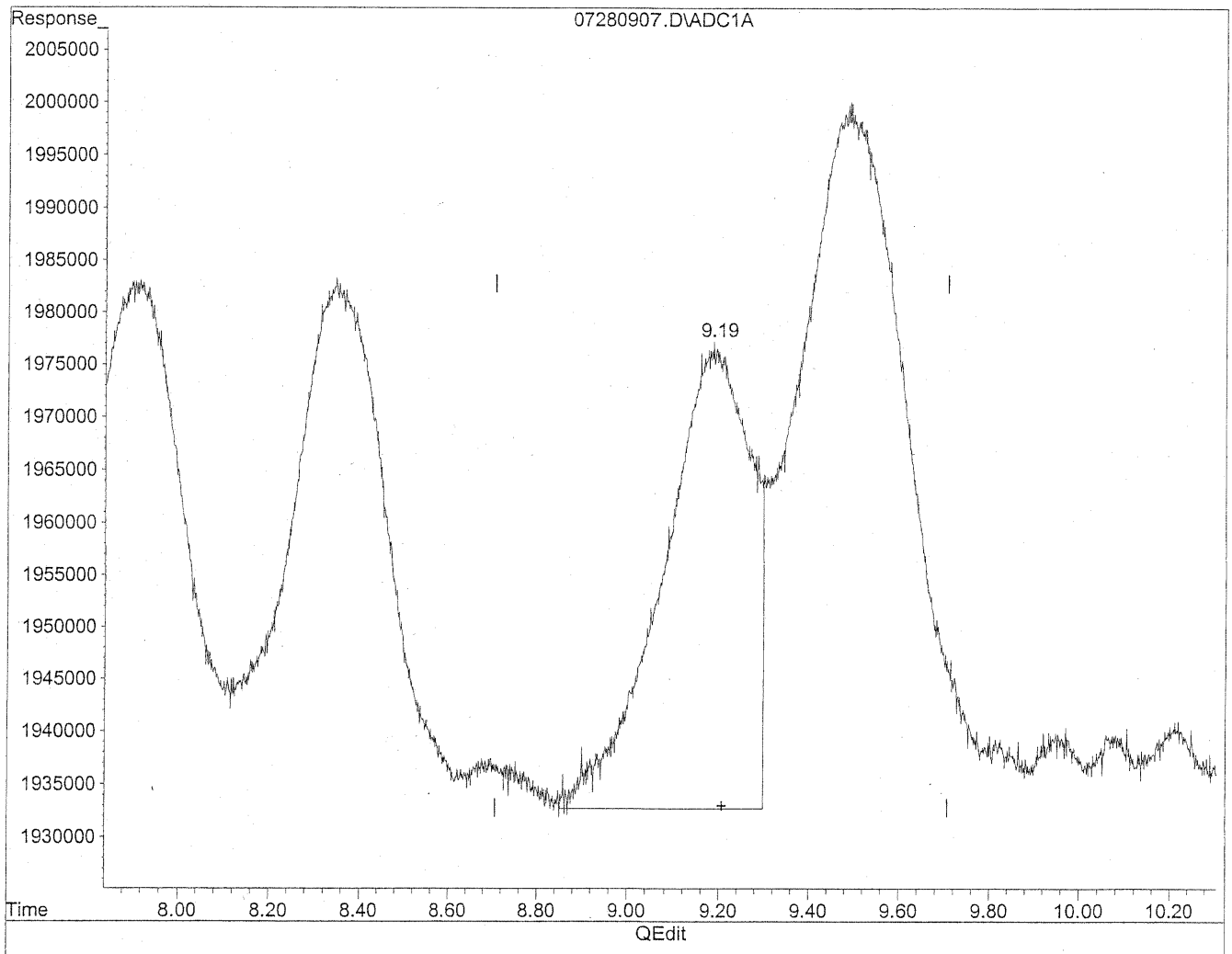
(+) = Expected Retention Time

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

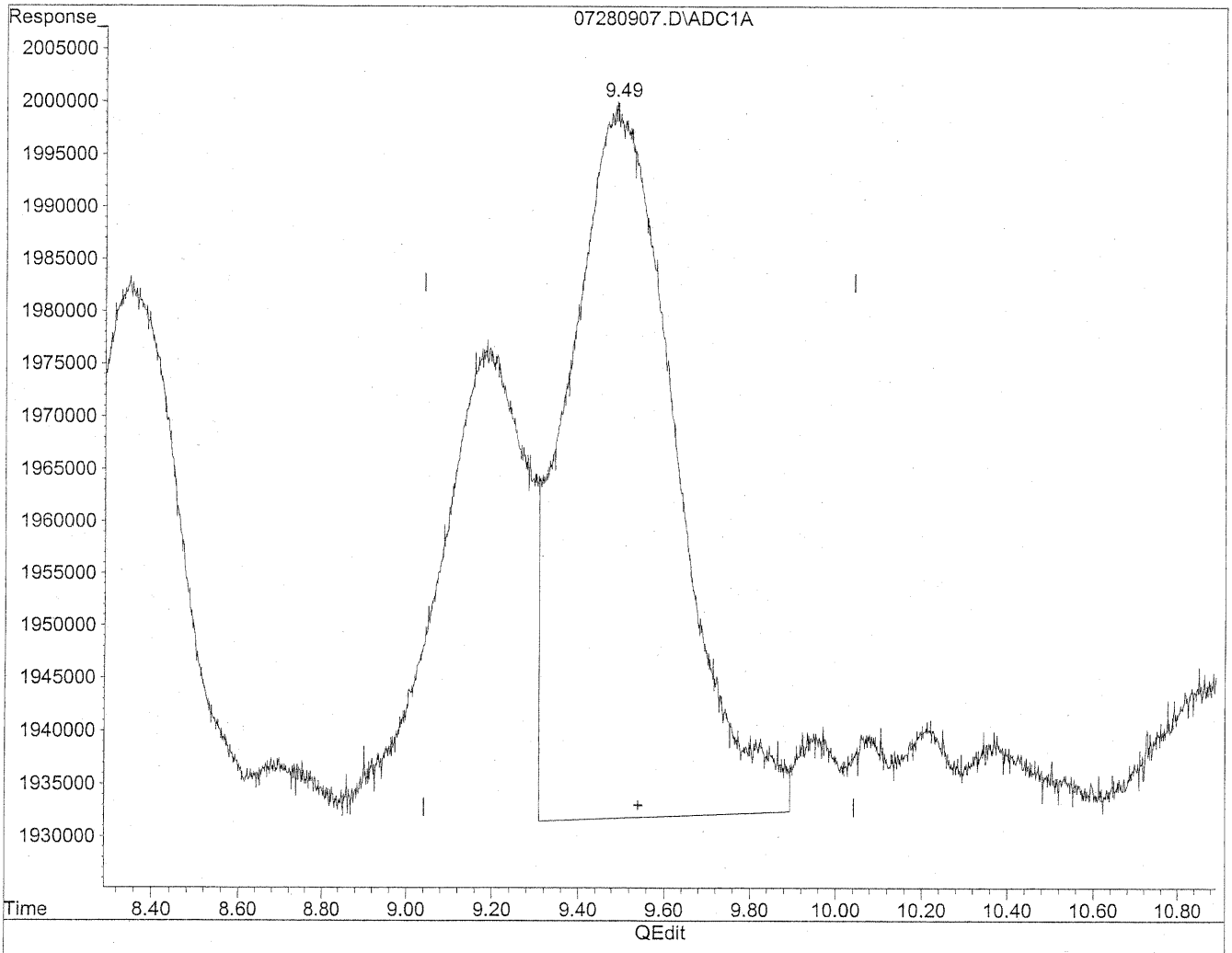
249

(+) = 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

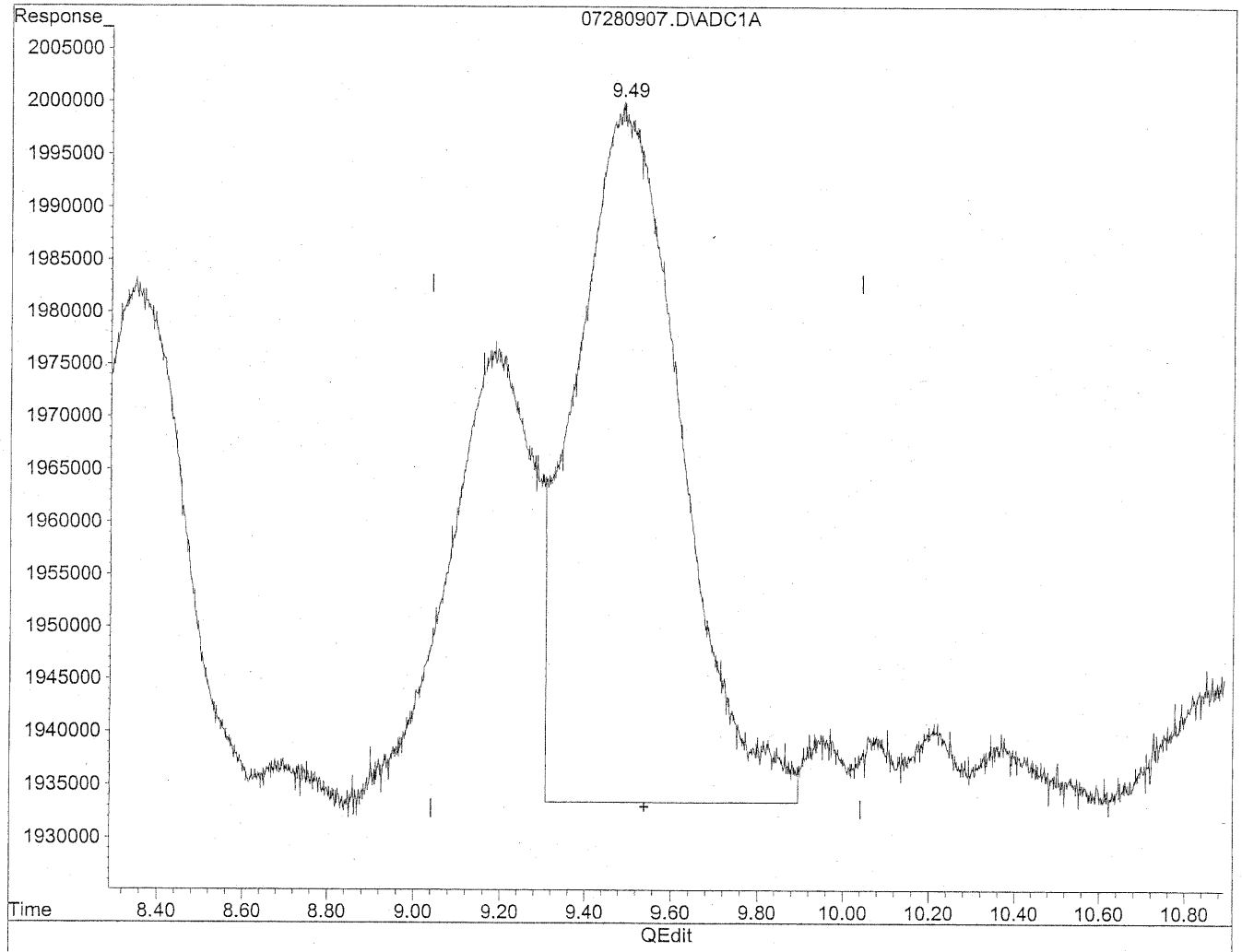
response 11738041

(+) = 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 BC

KR 7/29/09

251

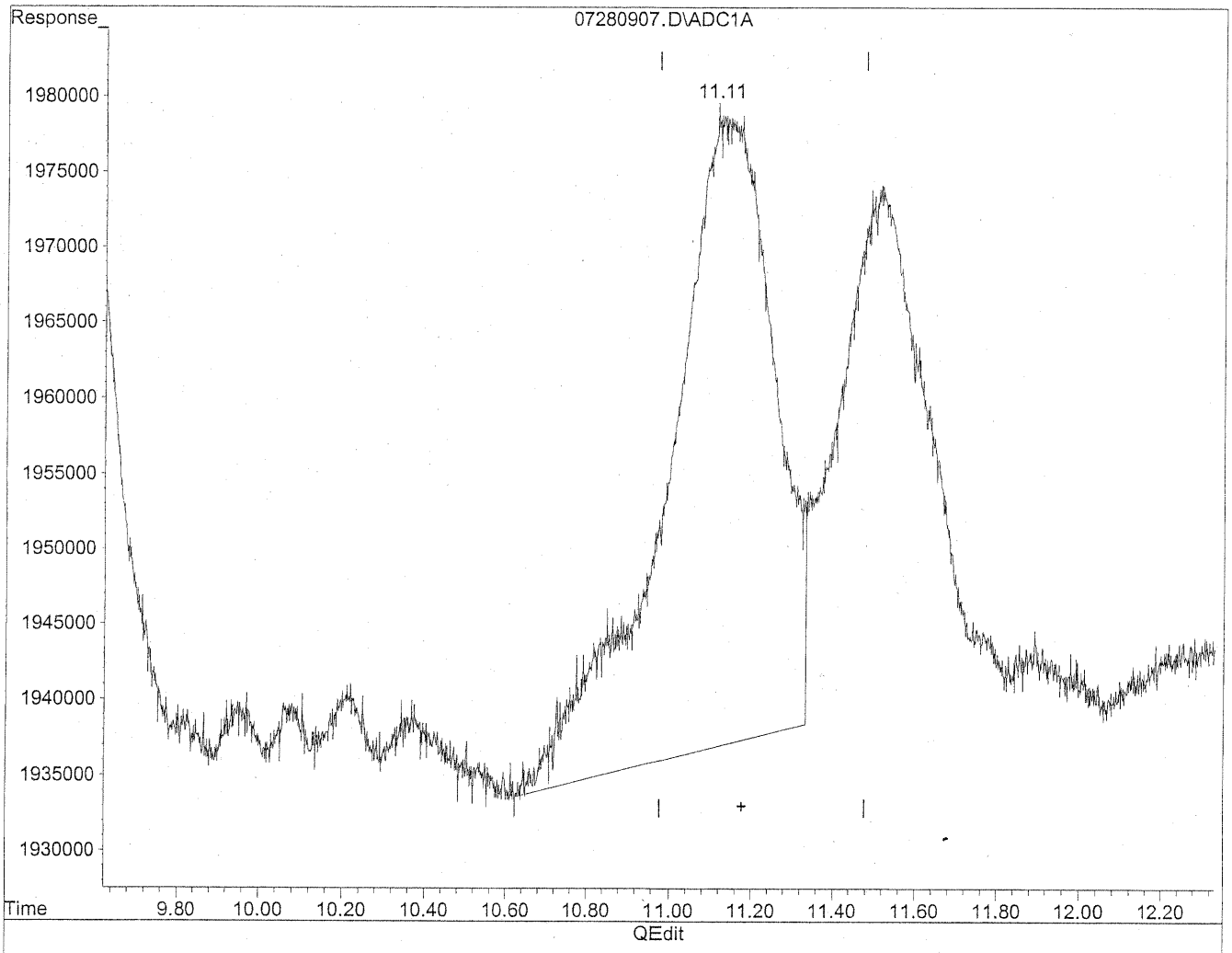
(+) = Expected Retention Time

07280907.D TO110709.M Tue Jul 28 14:32:53 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

11.14min 112.492ng/ml

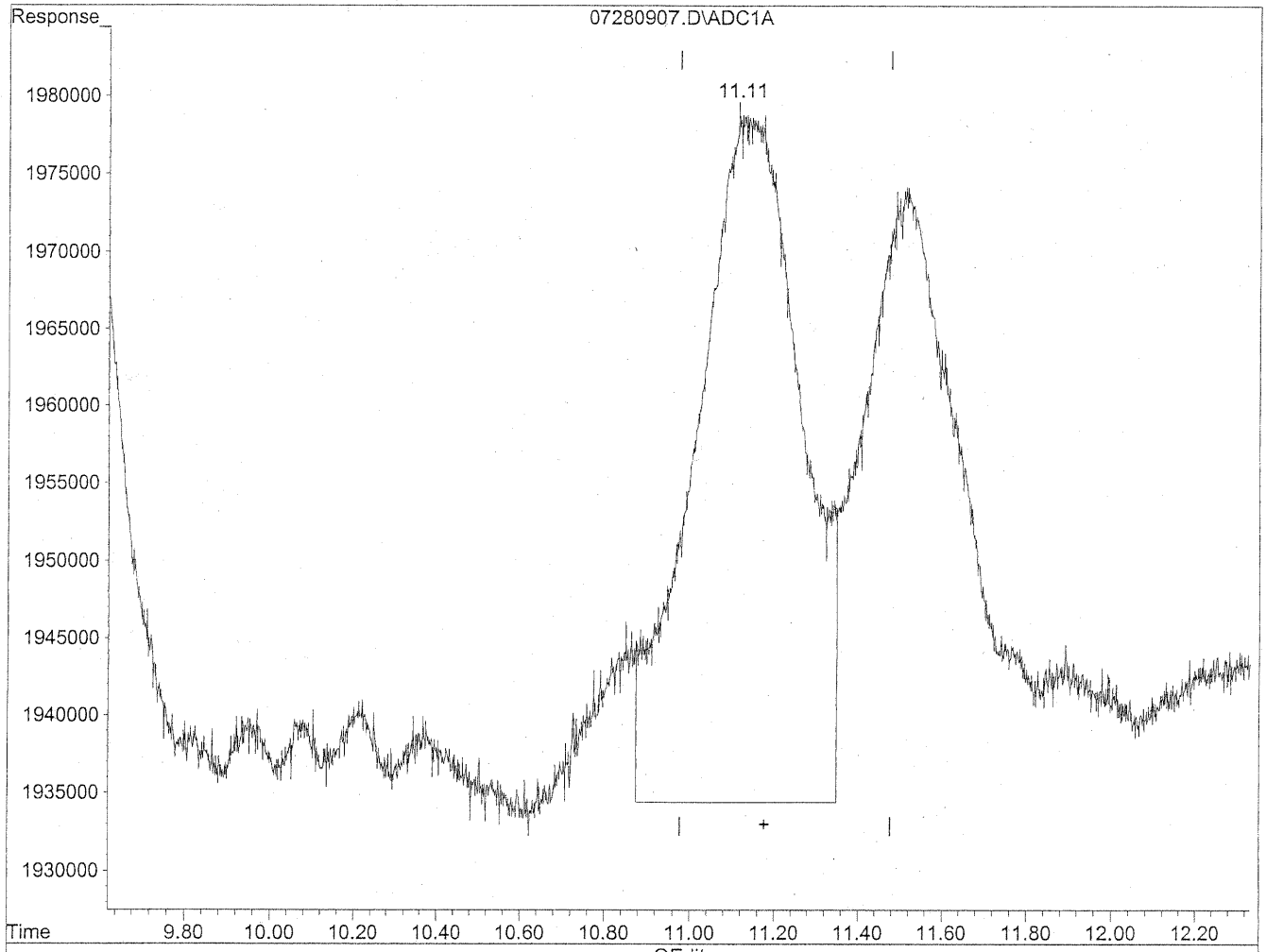
response 7552544

(+) = 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

11.11min 114.897ng/ml m

response 7714022

HC standard STL

KR 7/29/09

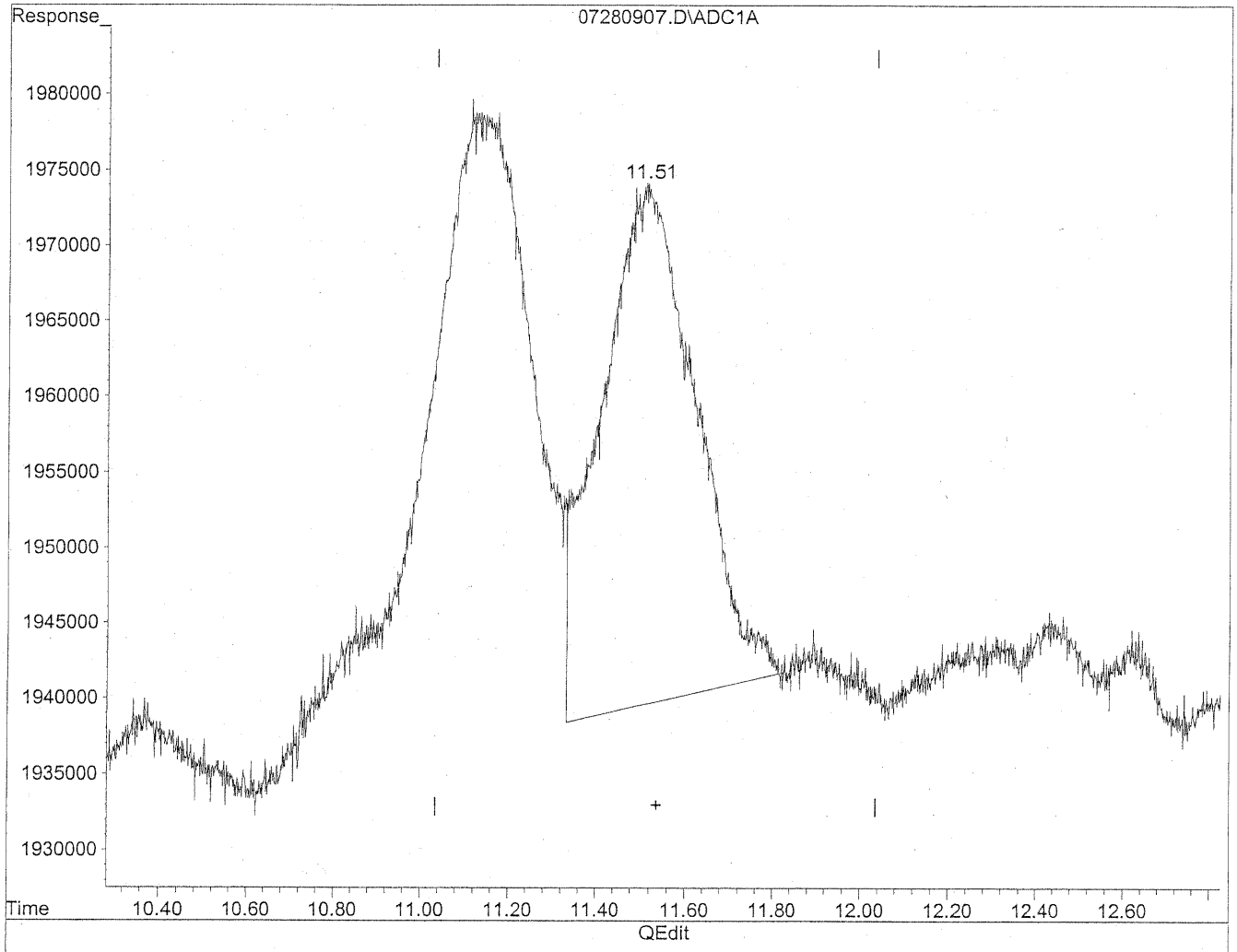
253

(+) = 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

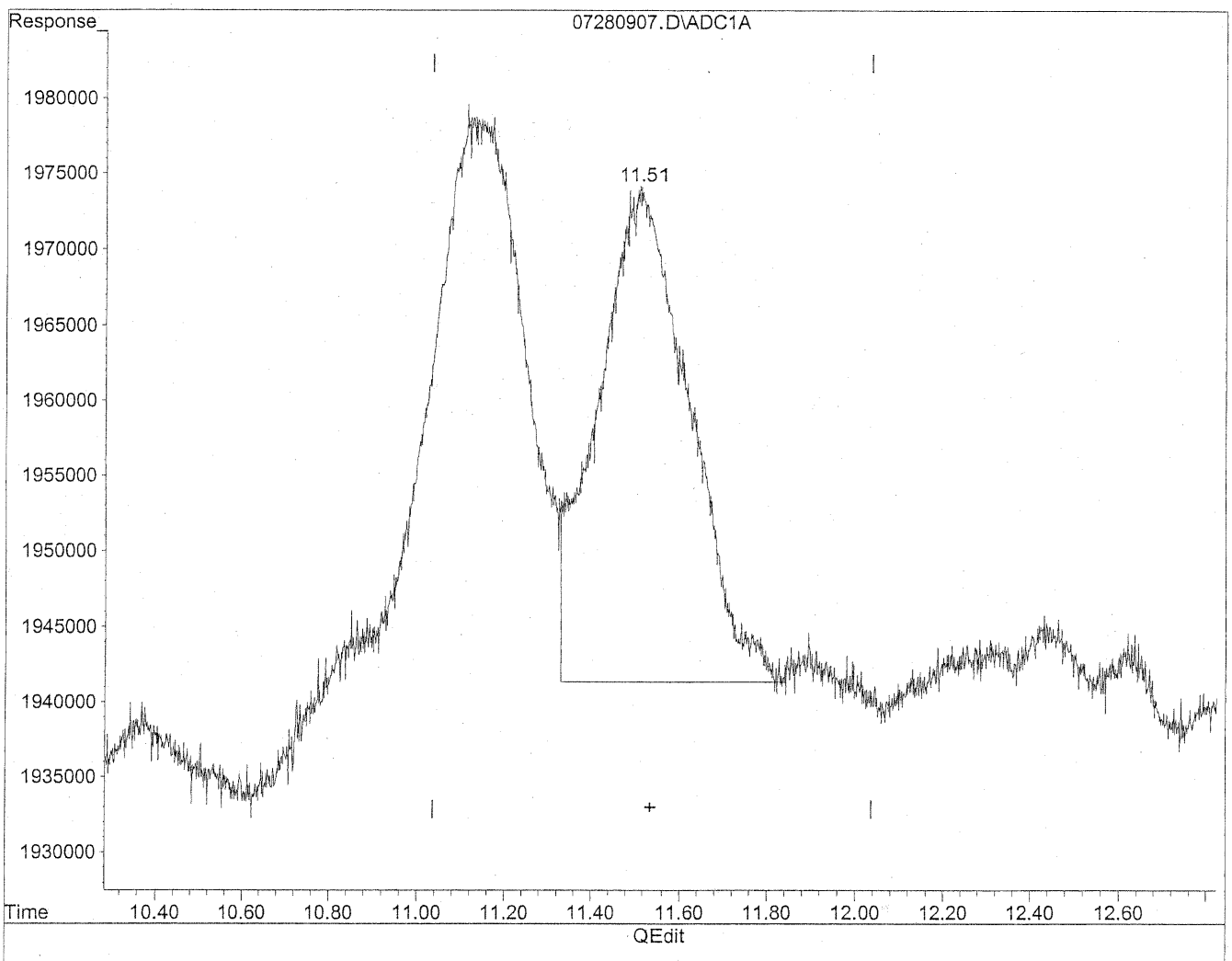
254

(+) = 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

11.51min 91.178ng/ml m

response 4735227

HL 7/28/09 BL

KR 7/29/09

255

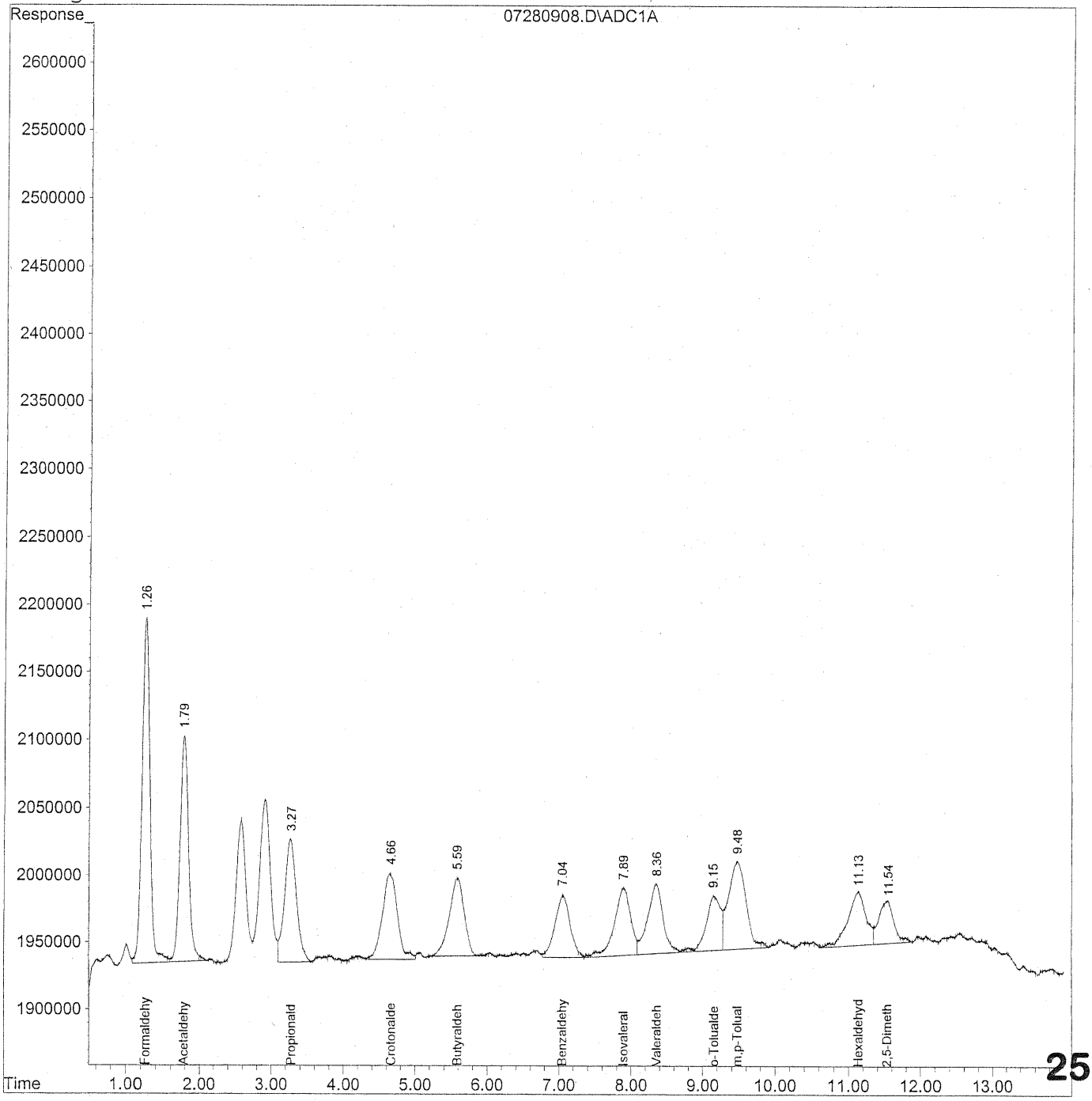
(+) = Expected Retention Time

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 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



256

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

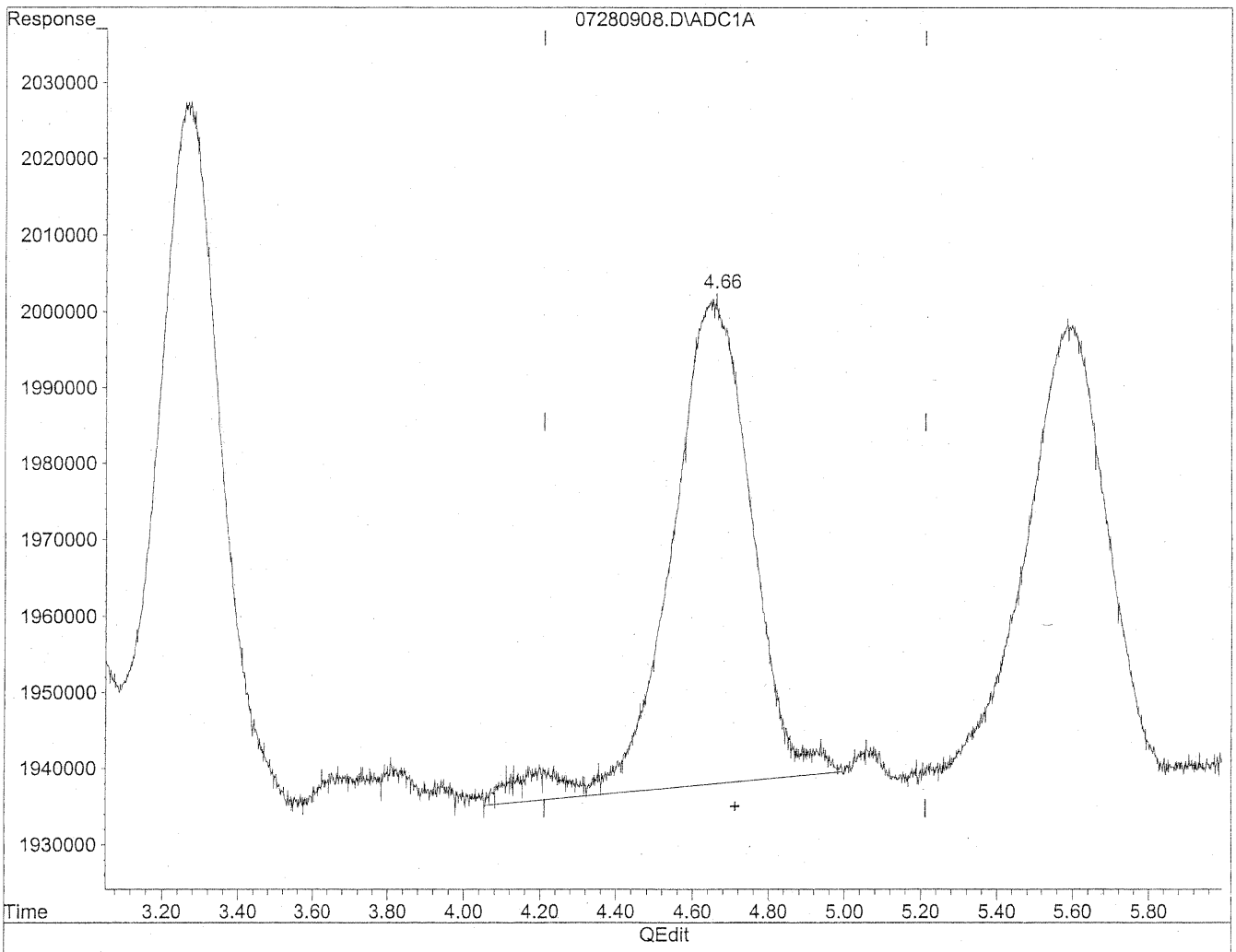
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.26	18400032	104.788	ng/ml
2)	Acetaldehyde	1.79	13737532	101.835	ng/ml
3)	Propionaldehyde	3.27	10633406	103.442	ng/ml
4)	Crotonaldehyde	4.66	9424529	85.247	ng/mlm
5)	Butyraldehyde	5.59	8463028	105.163	ng/ml
6)	Benzaldehyde	7.04	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

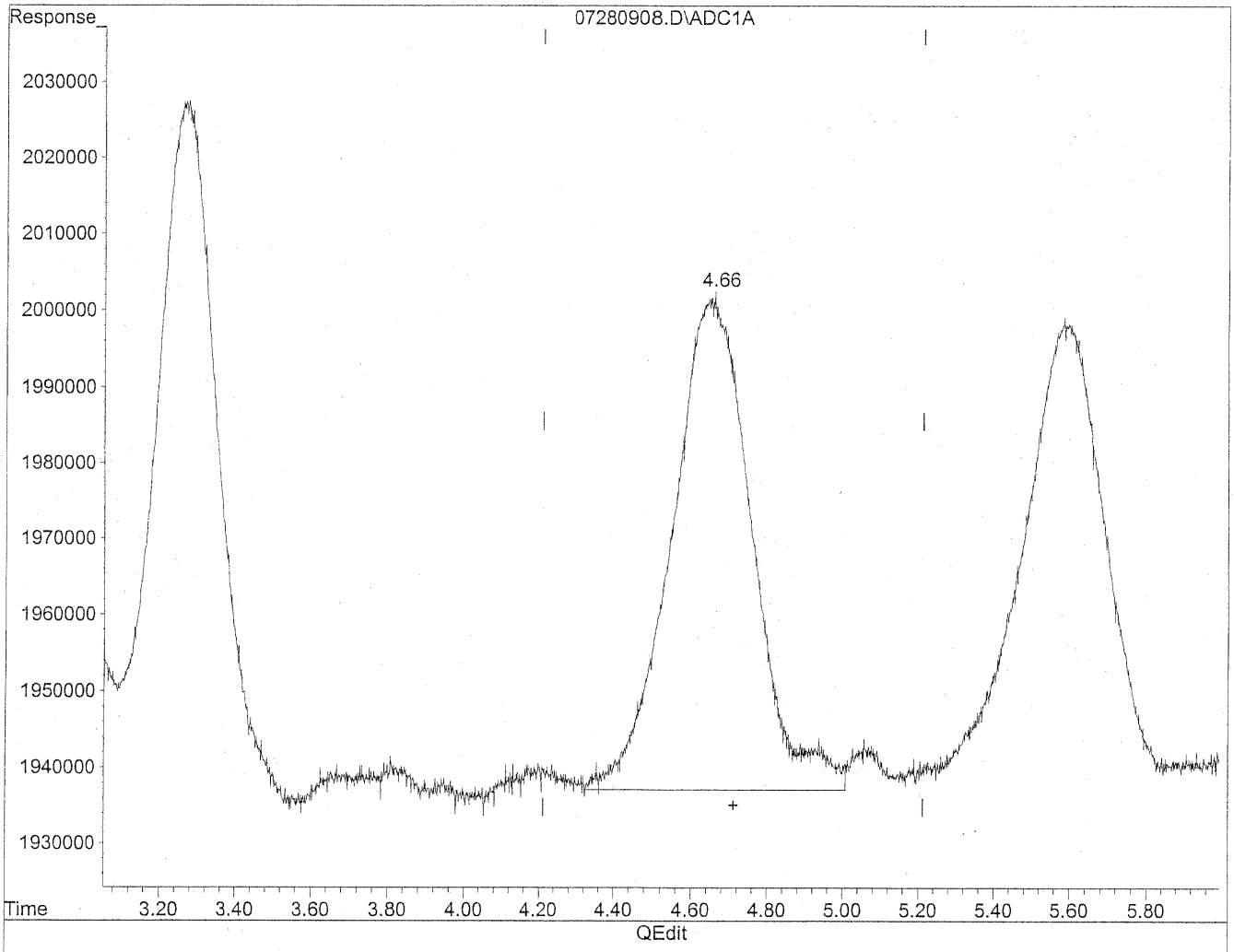
258

(+) = 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

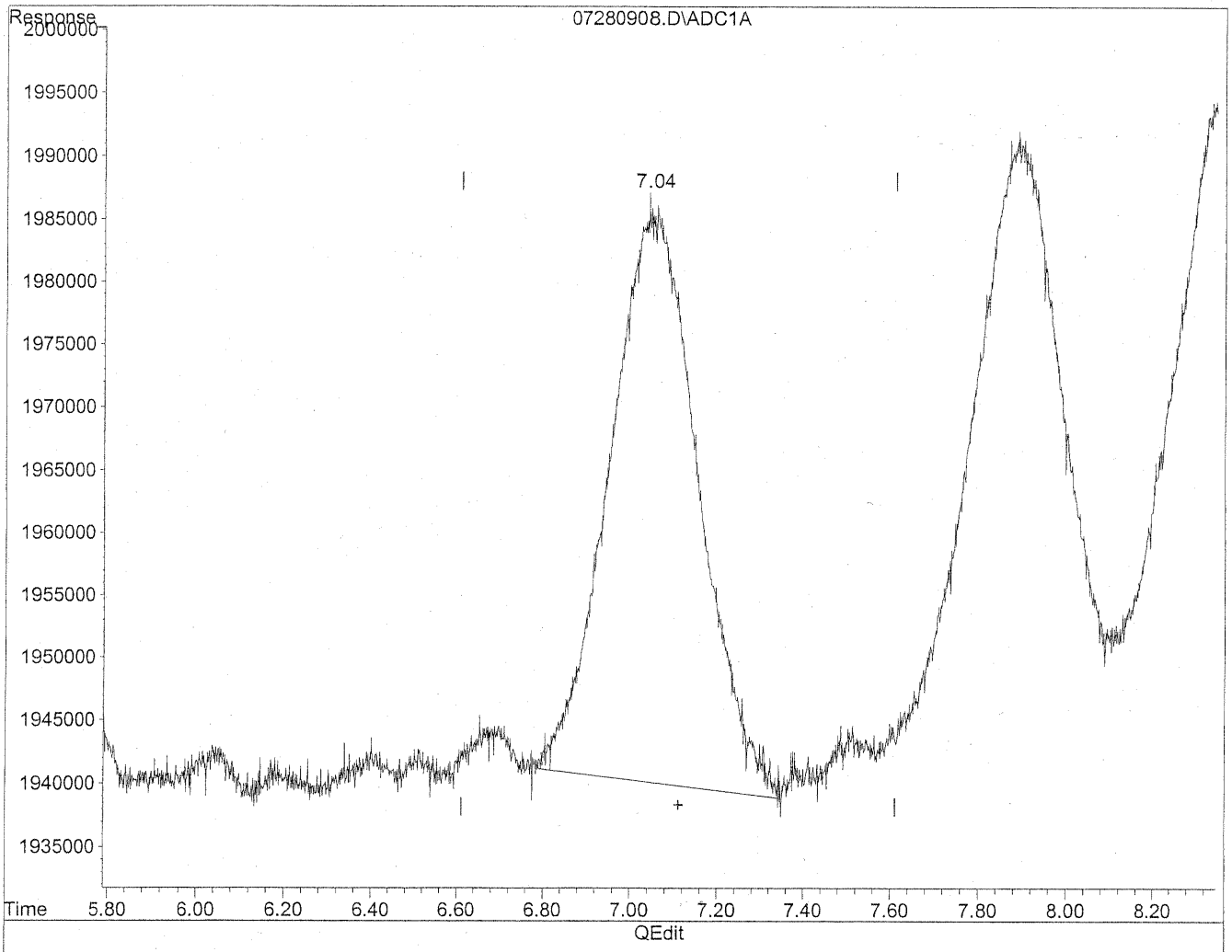
259

(+) = 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

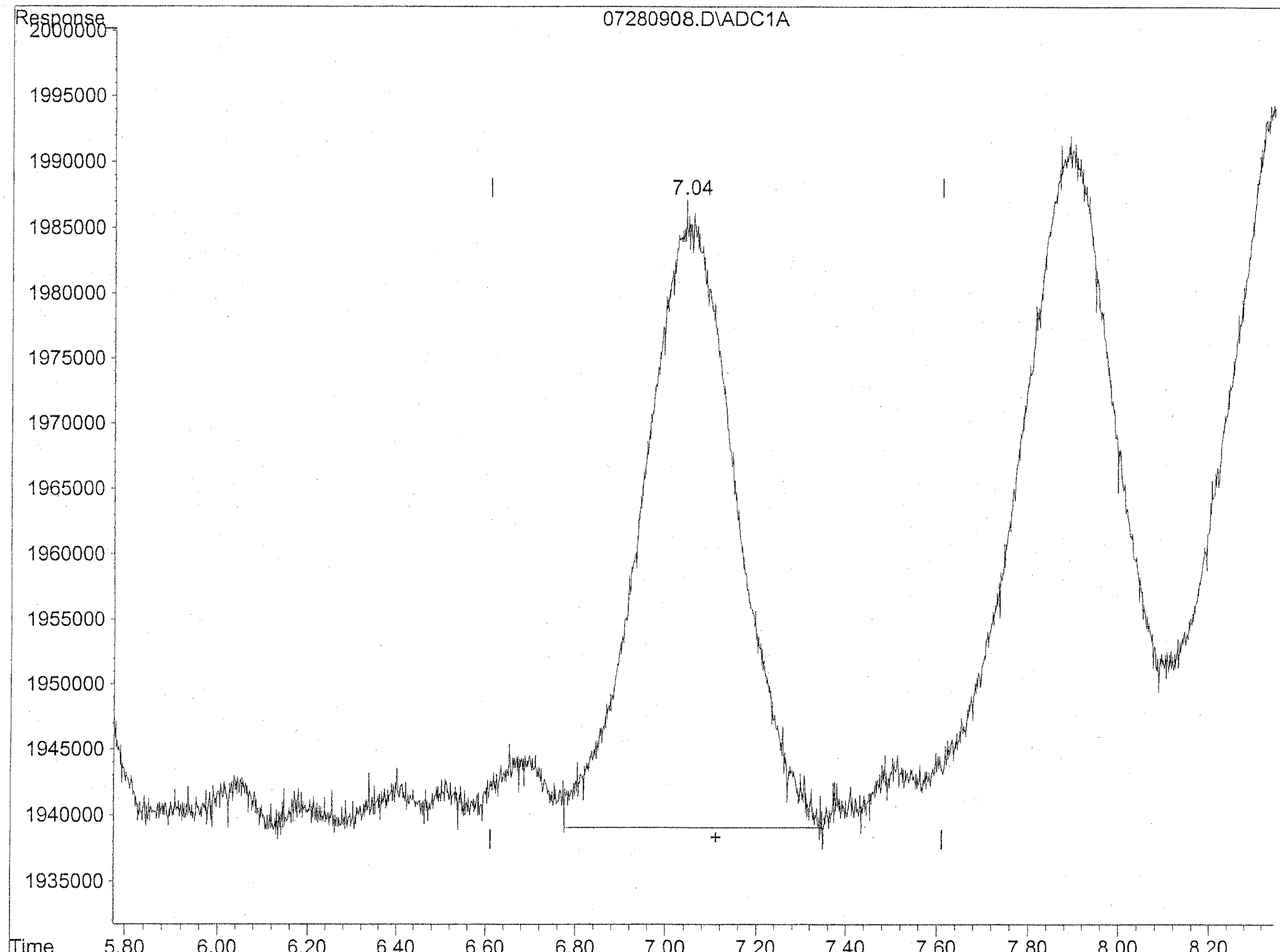
(+) = Expected Retention Time

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

KR 7/29/09

261

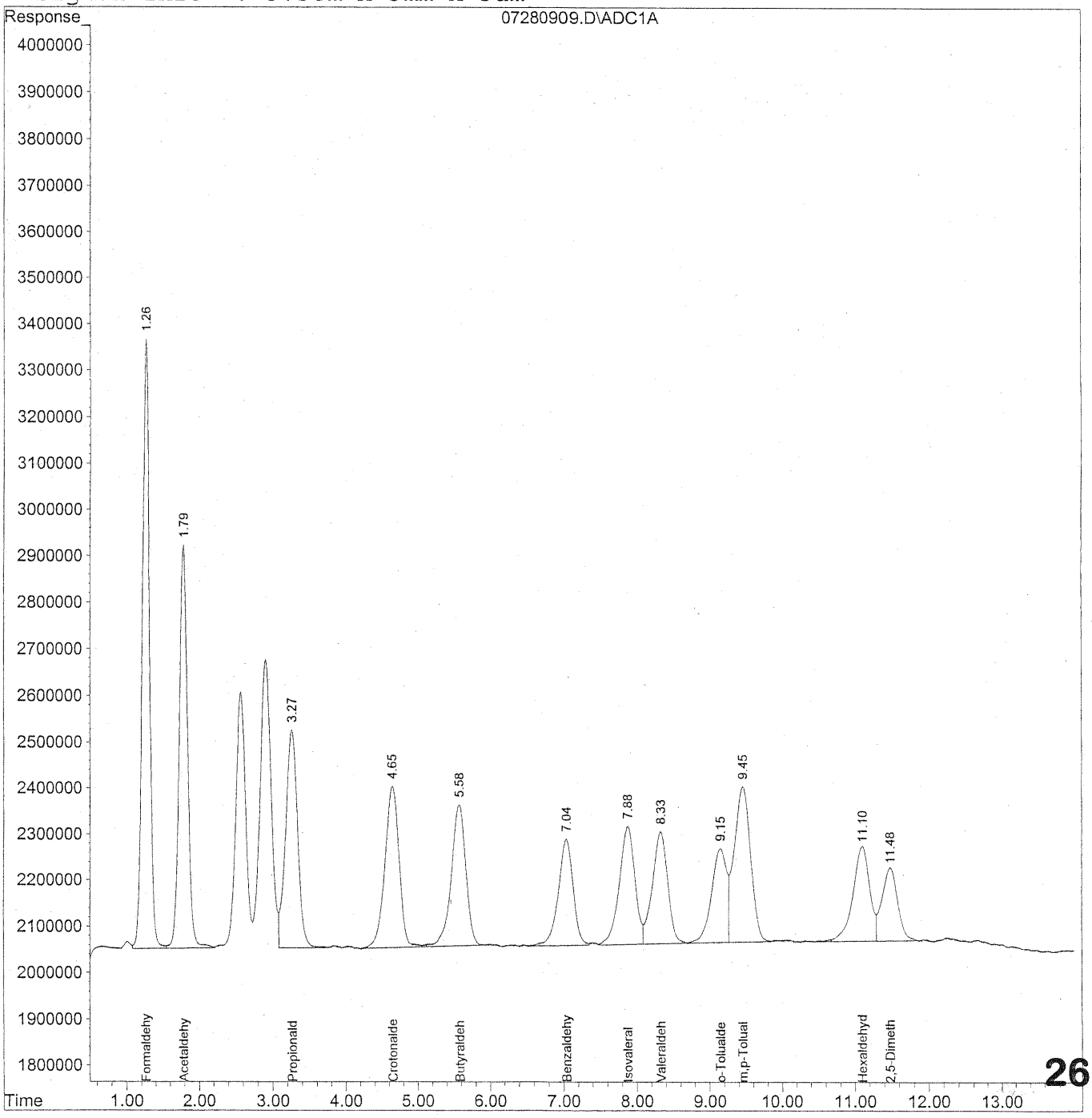
(+) = Expected Retention Time

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



262

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

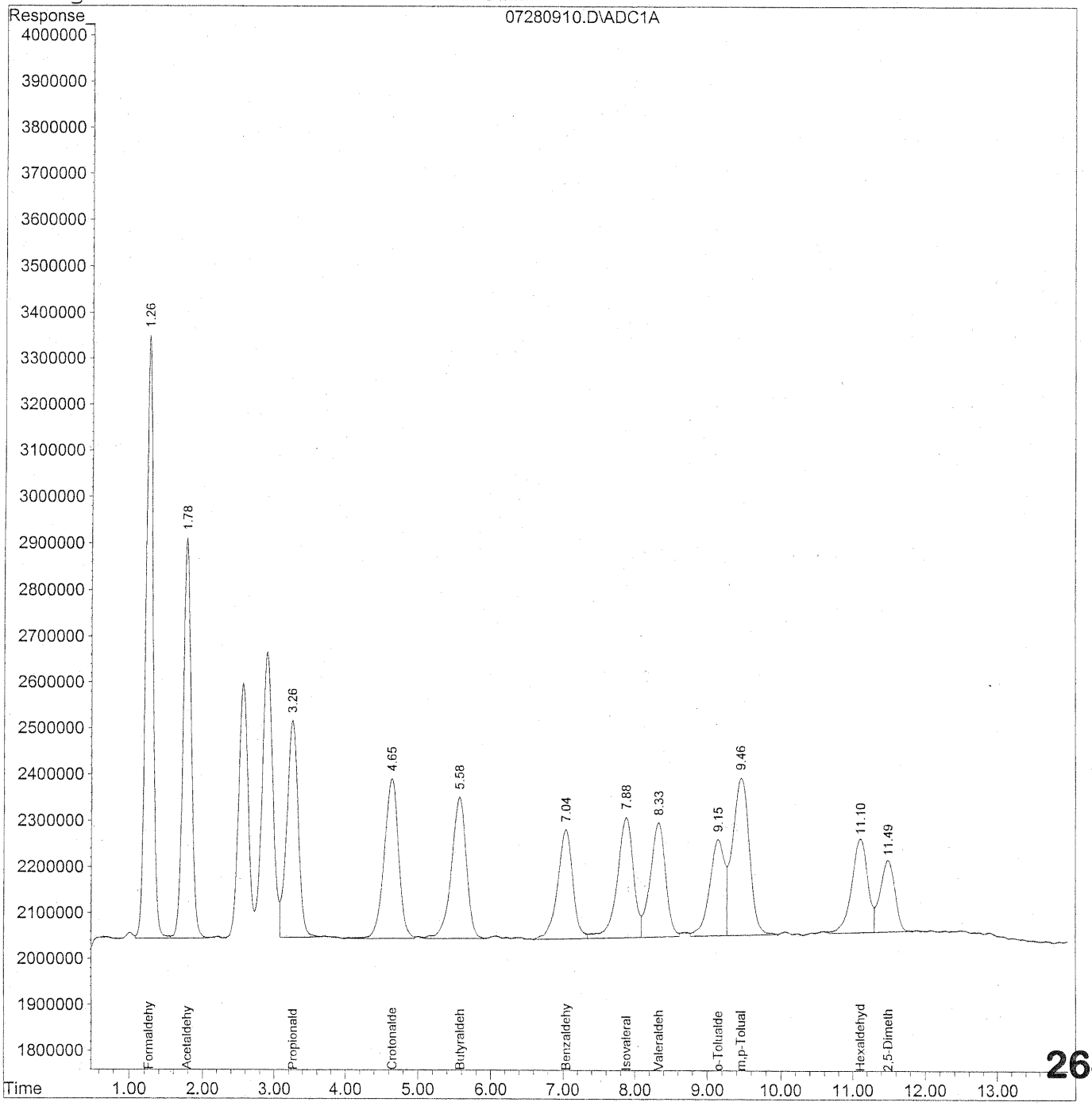
(f)=RT Delta > 1/2 Window (m)=manual int.

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



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

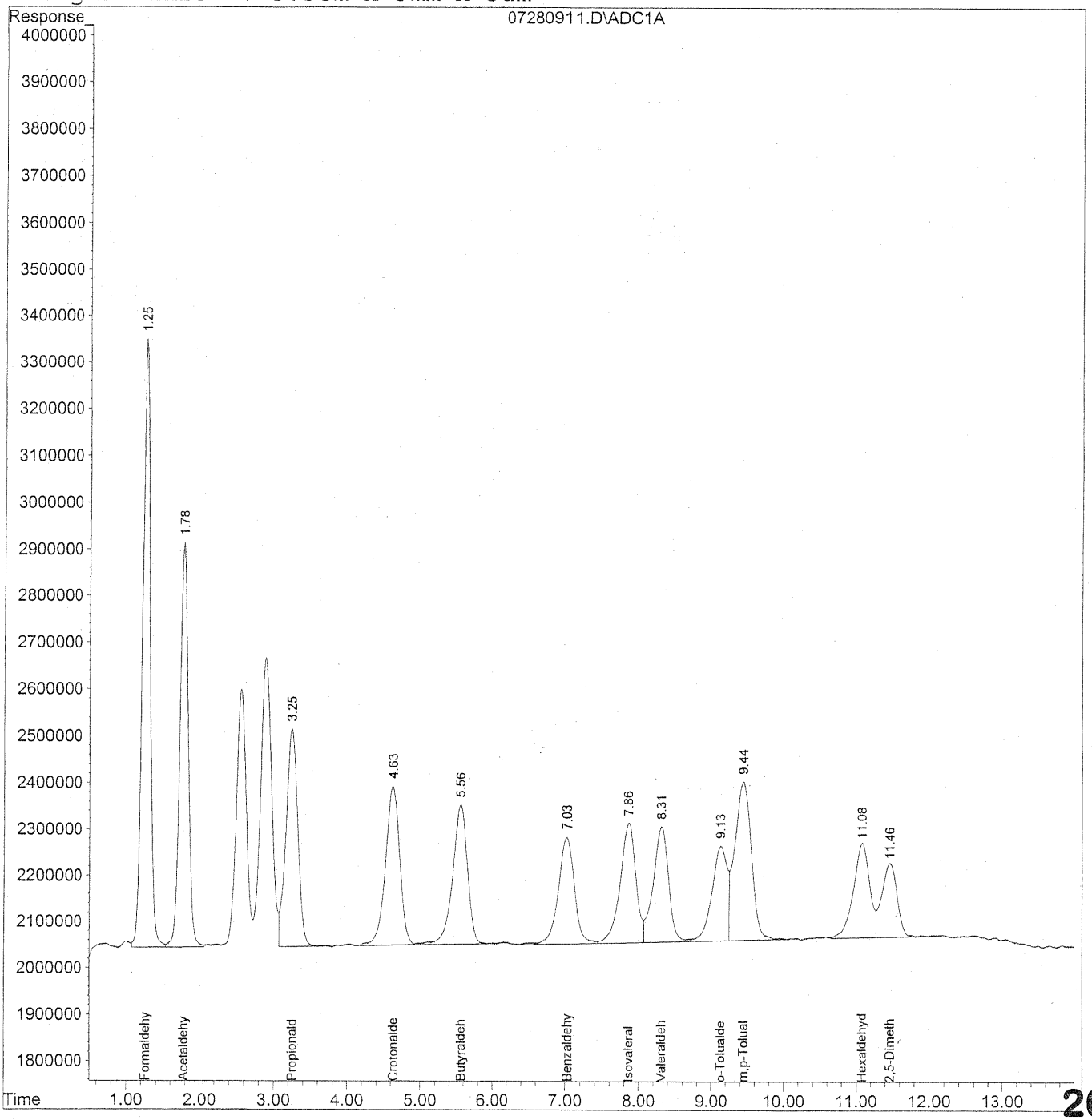
(f)=RT Delta > 1/2 Window (m)=manual int.

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



266

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

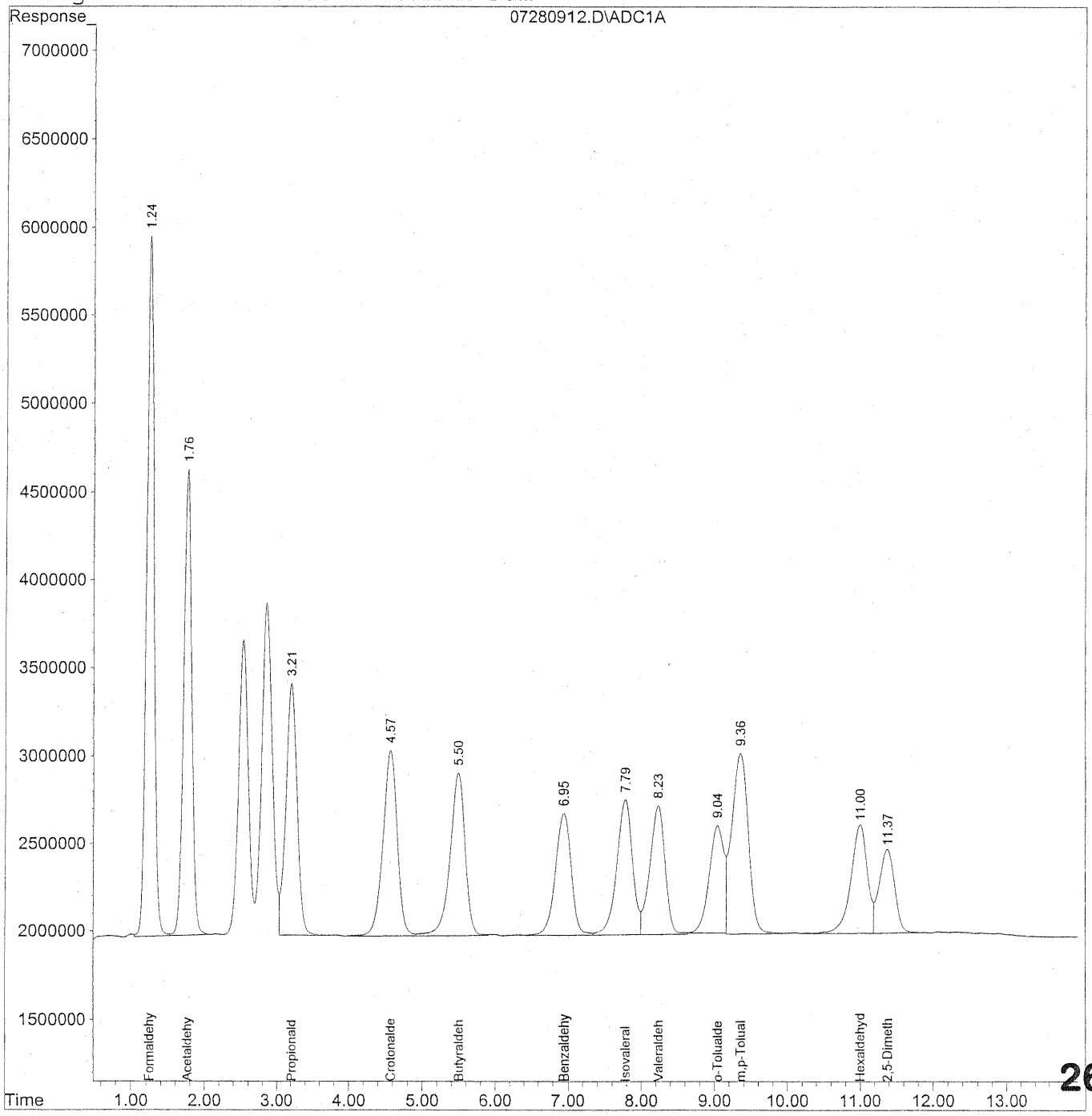
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

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



268

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

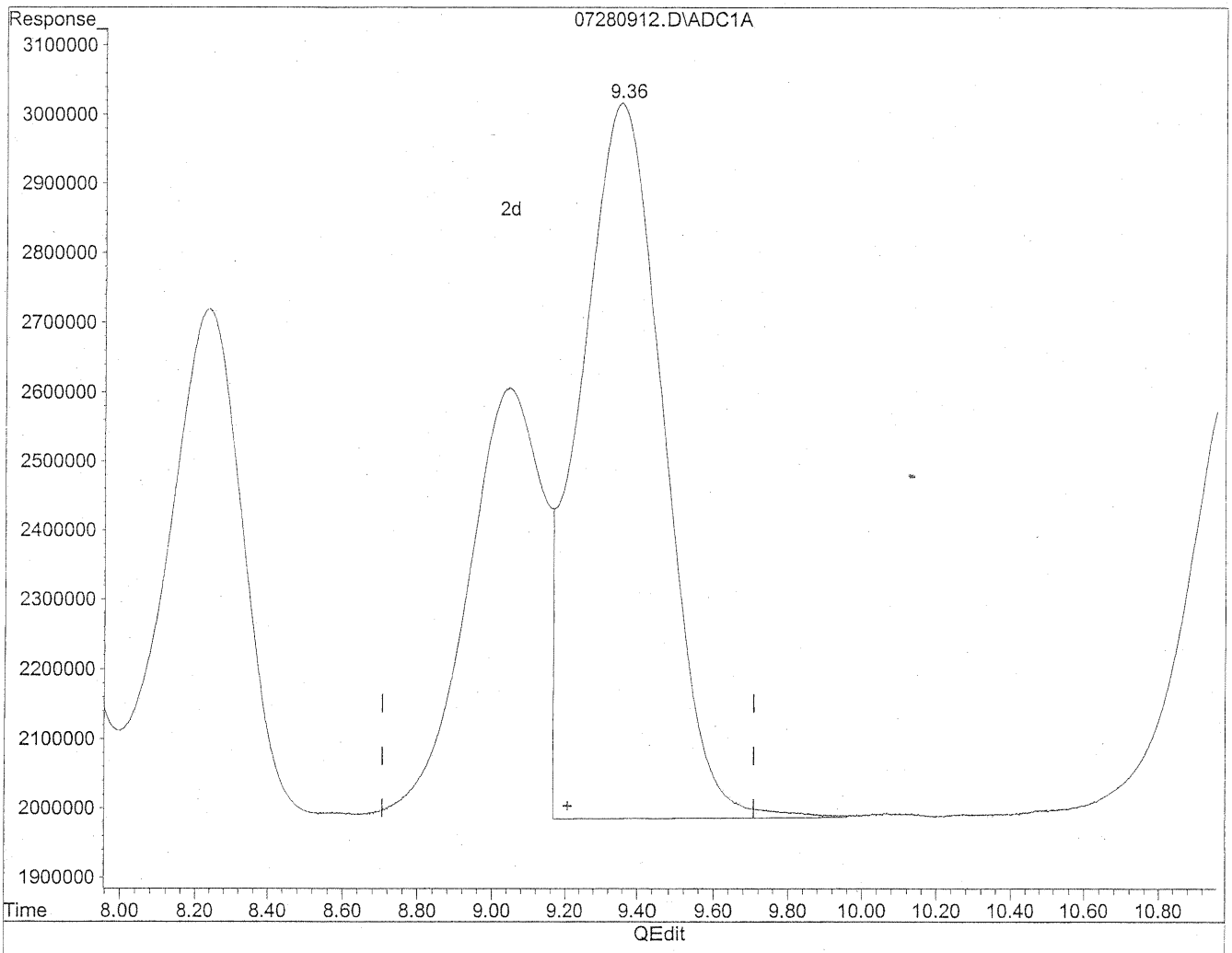
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.24	275380897	1568.292	ng/ml
2)	Acetaldehyde	1.76	209374751	1552.082	ng/ml
3)	Propionaldehyde	3.21	159030091	1547.054	ng/ml
4)	Crotonaldehyde	4.57	143227783	1295.530	ng/ml
5)	Butyraldehyde	5.50	134132687	1666.757	ng/ml
6)	Benzaldehyde	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

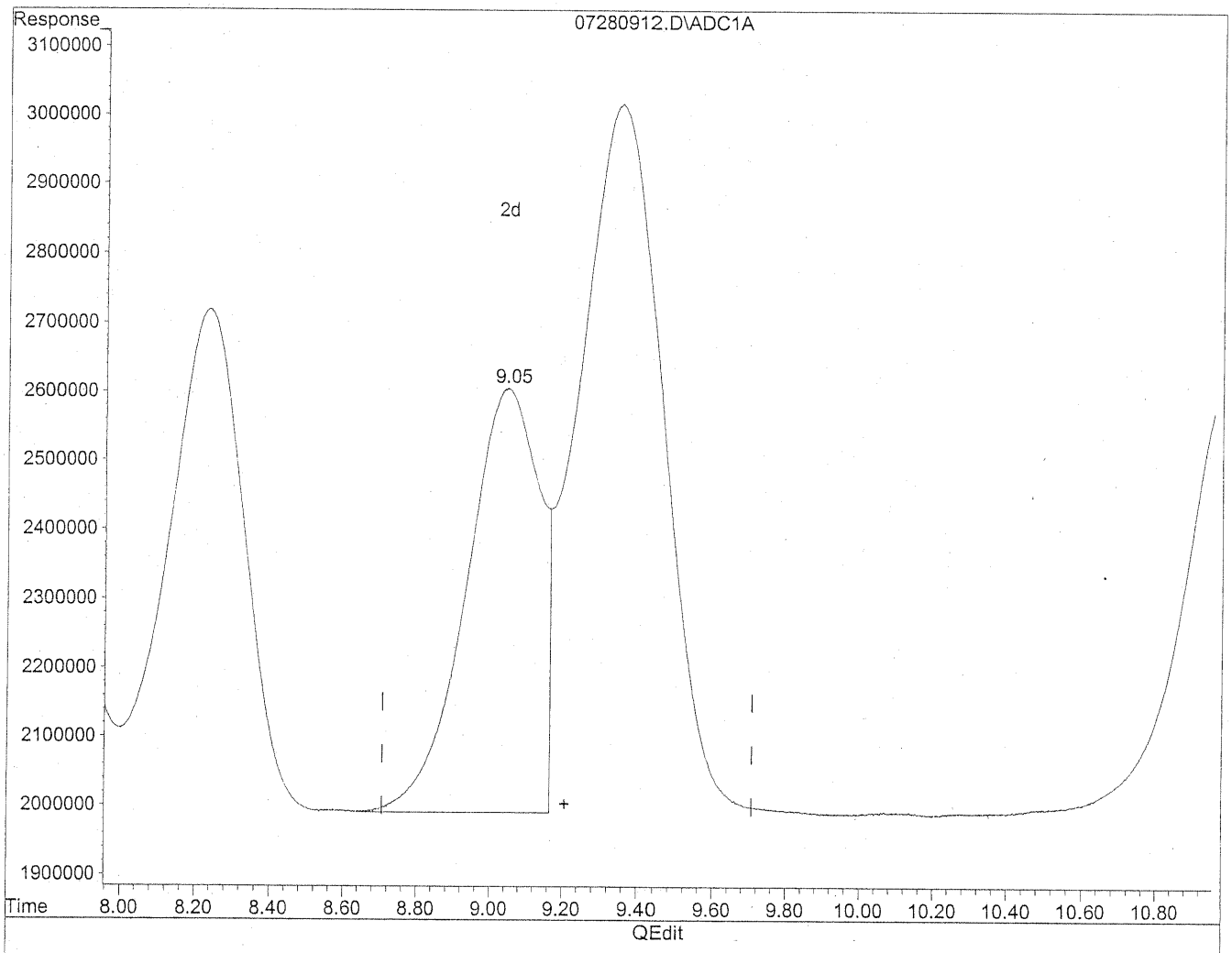
270

(+) = 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 WP

7/29/09

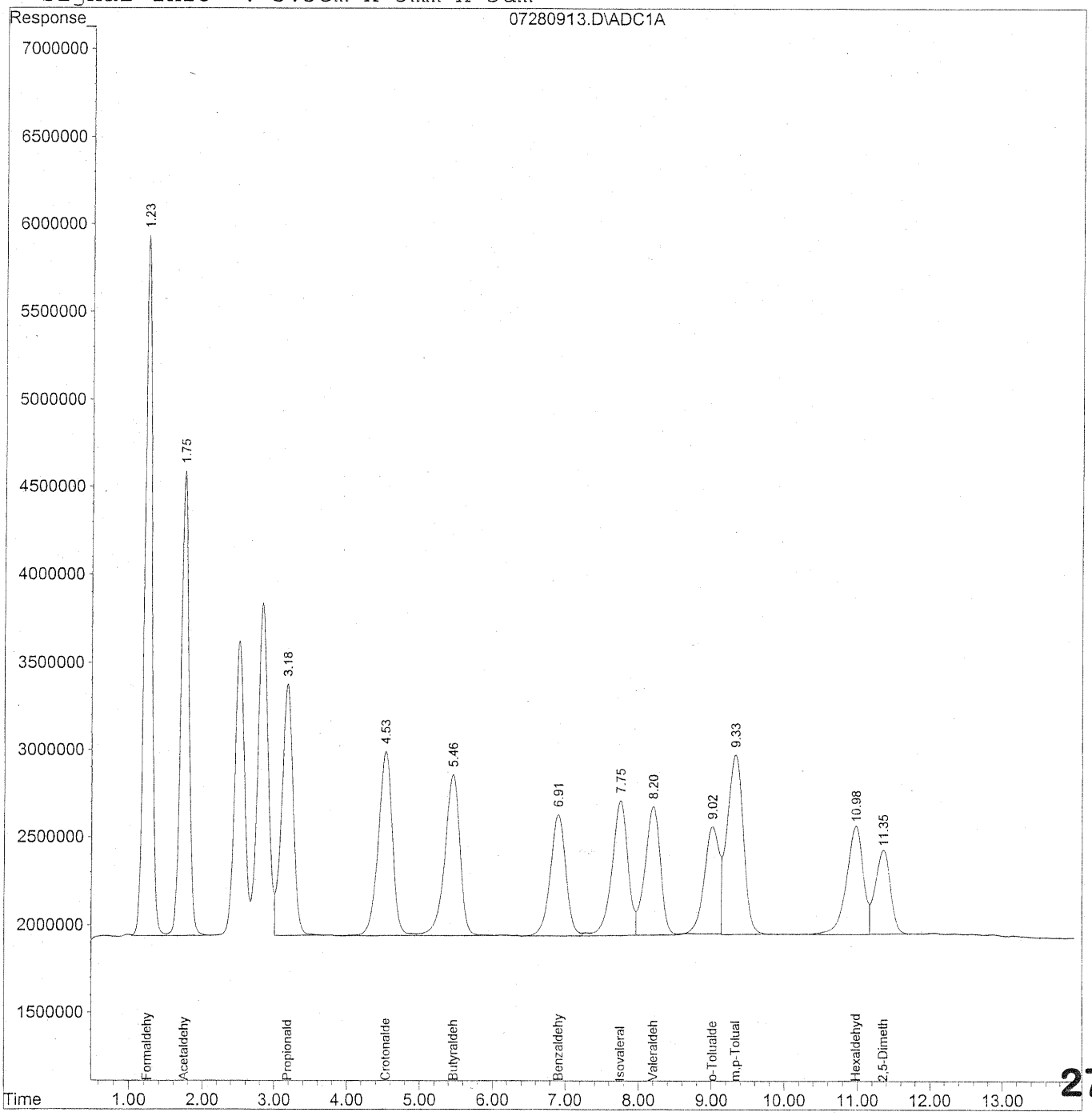
271

(+) = 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



272

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

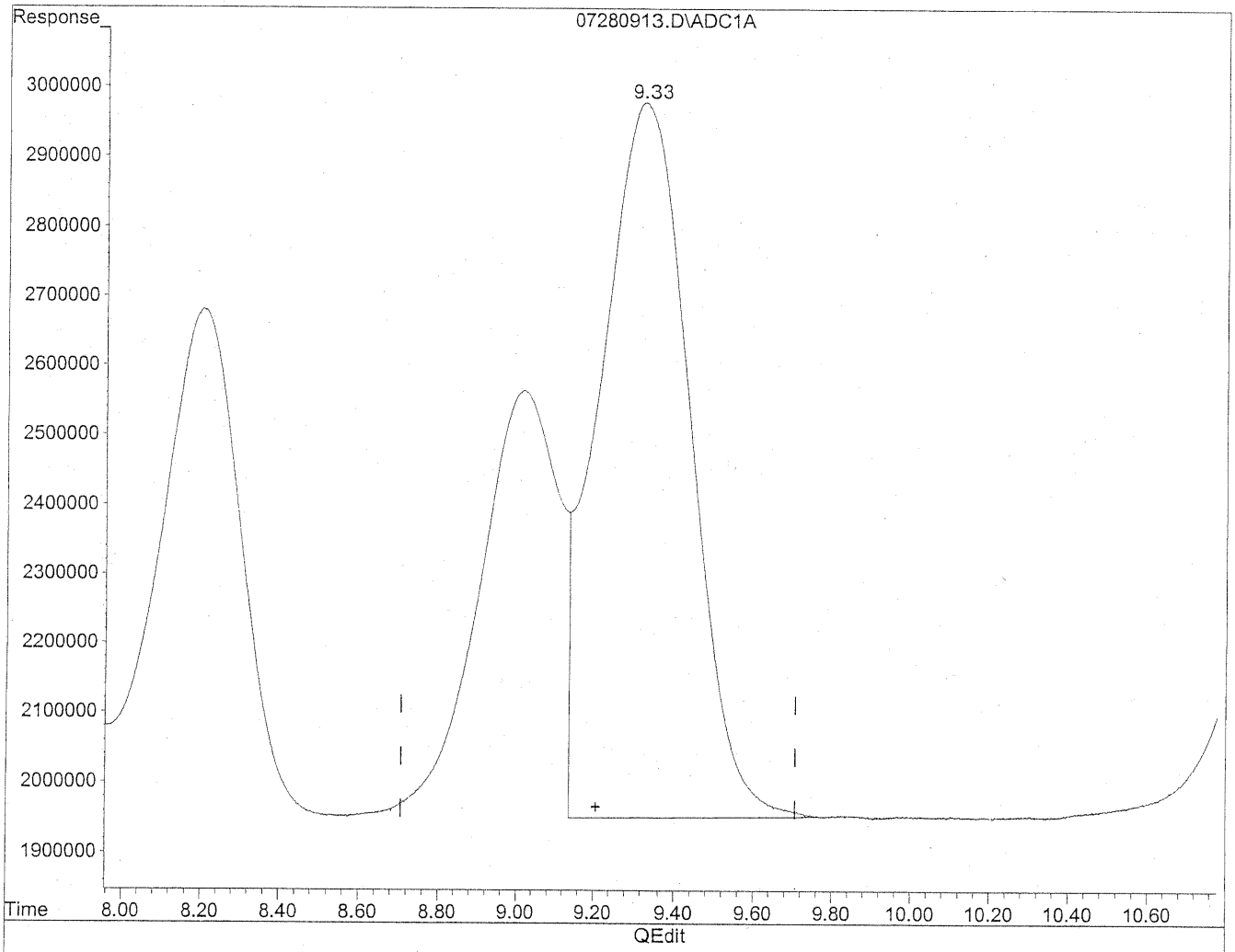
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.23	274724982	1564.557	ng/ml
2)	Acetaldehyde	1.75	209301649	1551.540	ng/ml
3)	Propionaldehyde	3.18	158919579	1545.979	ng/ml
4)	Crotonaldehyde	4.53	142112419	1285.442	ng/ml
5)	Butyraldehyde	5.46	132549734	1647.087	ng/ml
6)	Benzaldehyde	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

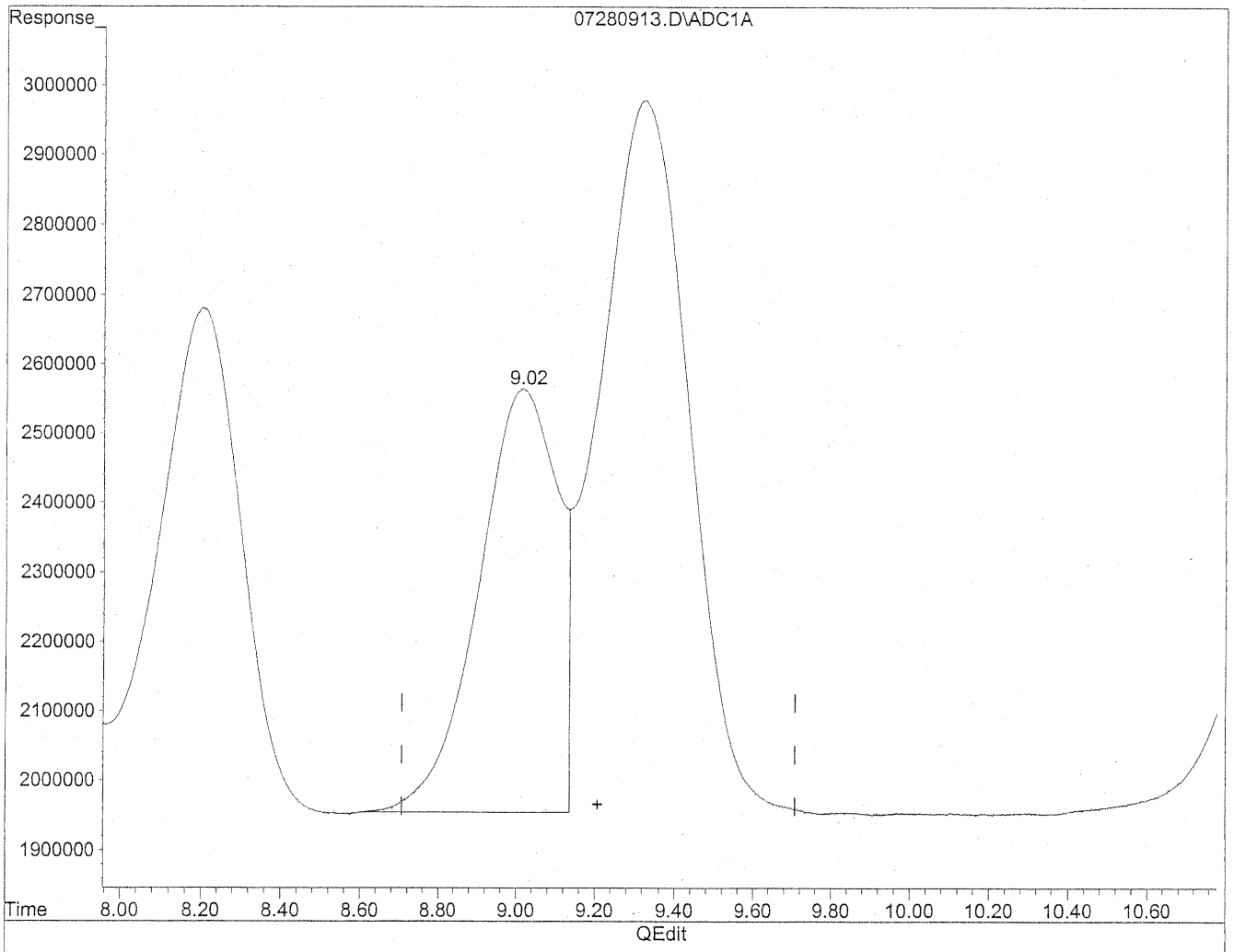
274

(+) = 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



(9) o-Tolualdehyde

9.02min 1595.318ng/ml m

response 85940120

7/28/09

7/29/09

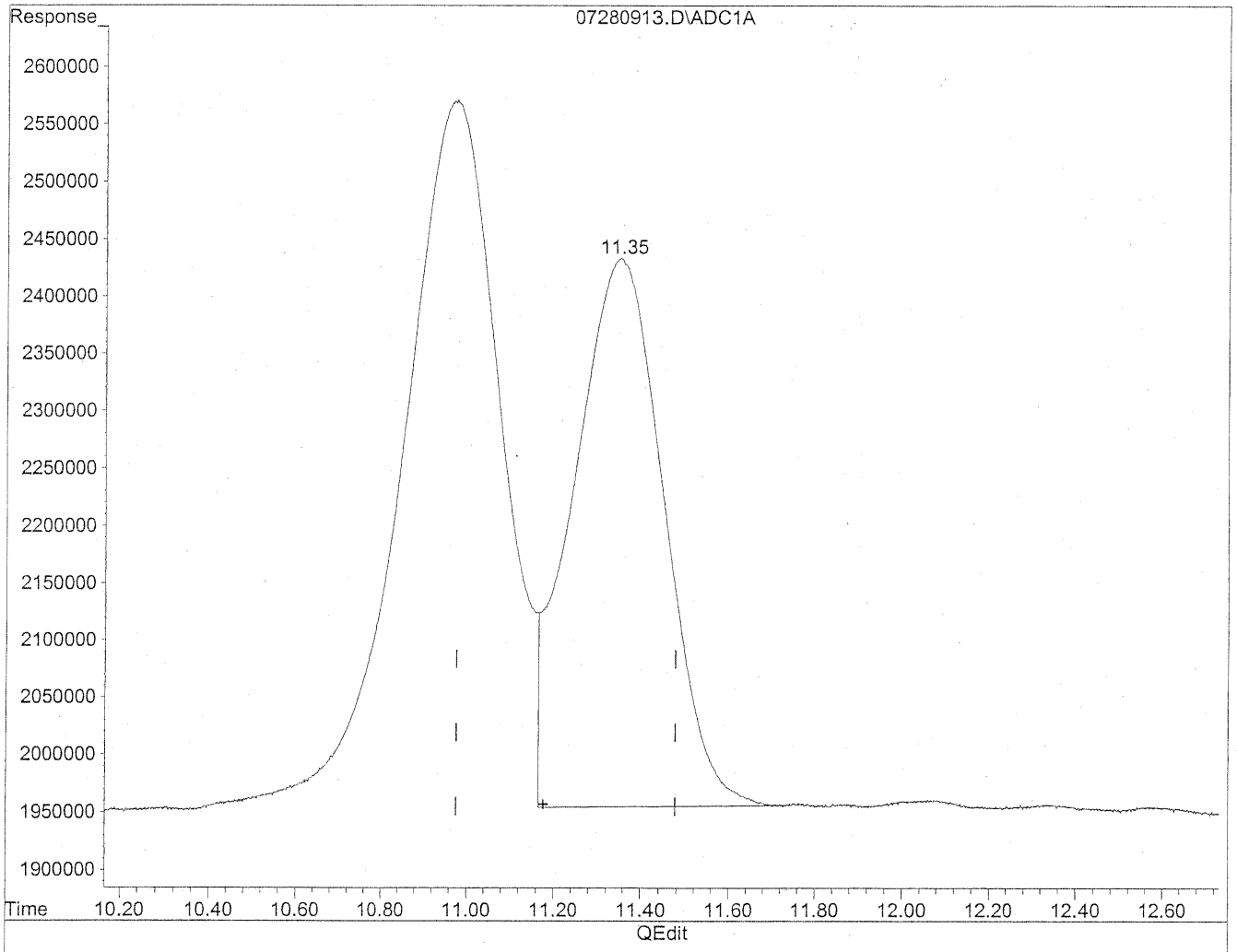
275

(+) = 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



(11) Hexaldehyde

11.35min 1025.842ng/ml

response 68873541

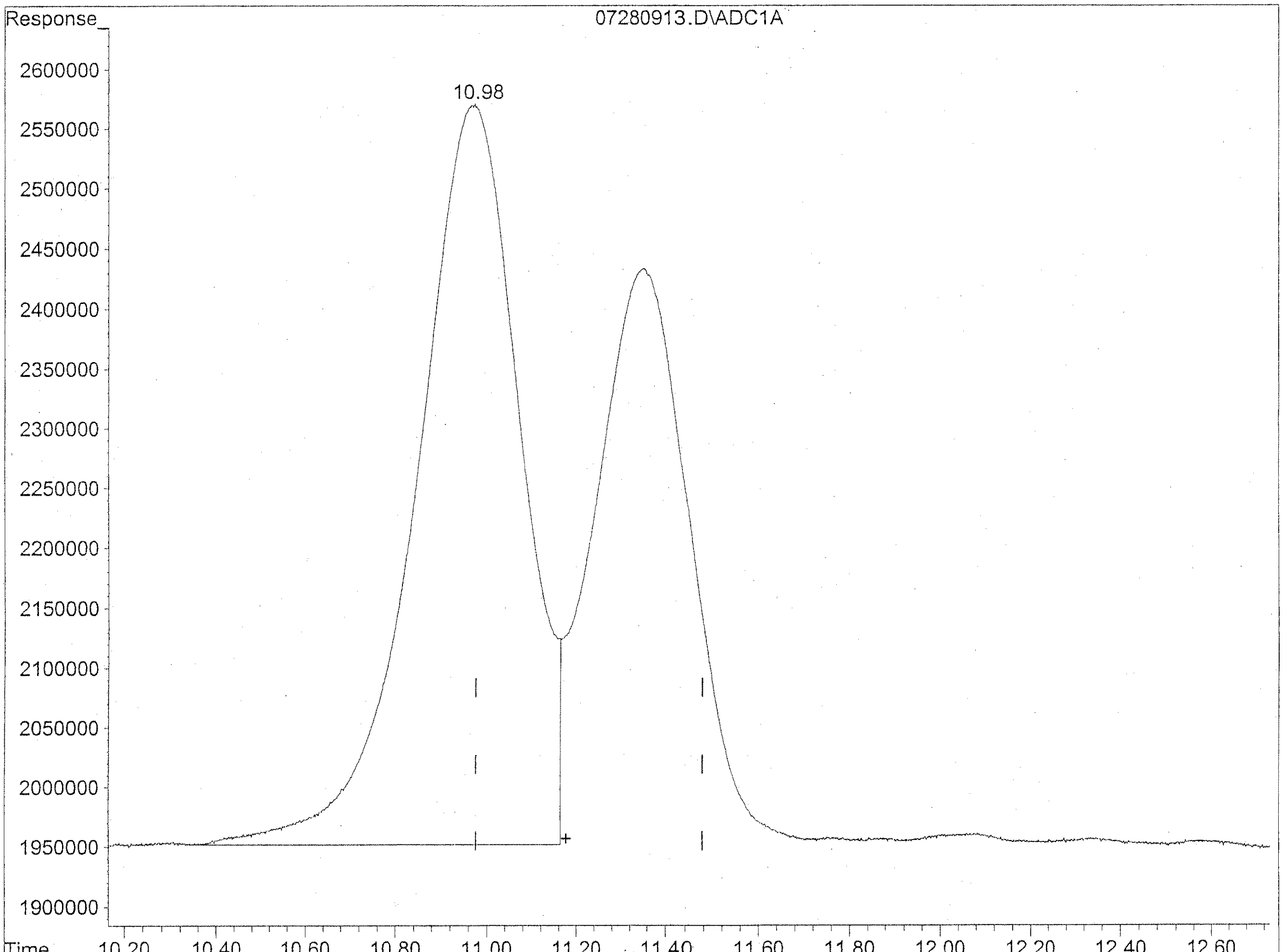
276

(+) = 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

HC 7/28/09 NR

KR 7/29/09

(+) = Expected Retention Time

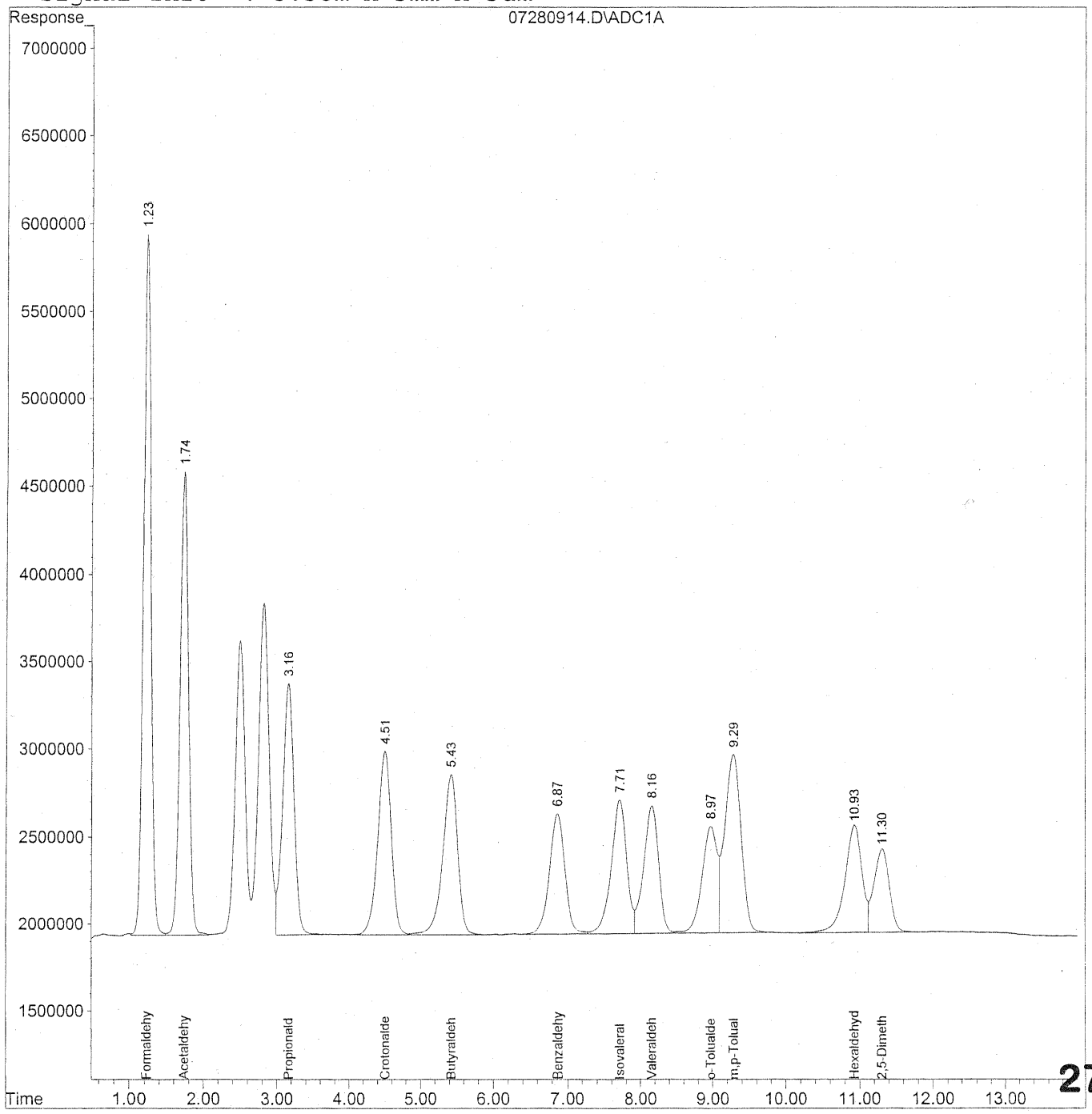
07280913.D   TO110709.M   Tue Jul 28 14:47:35 2009

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



278

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

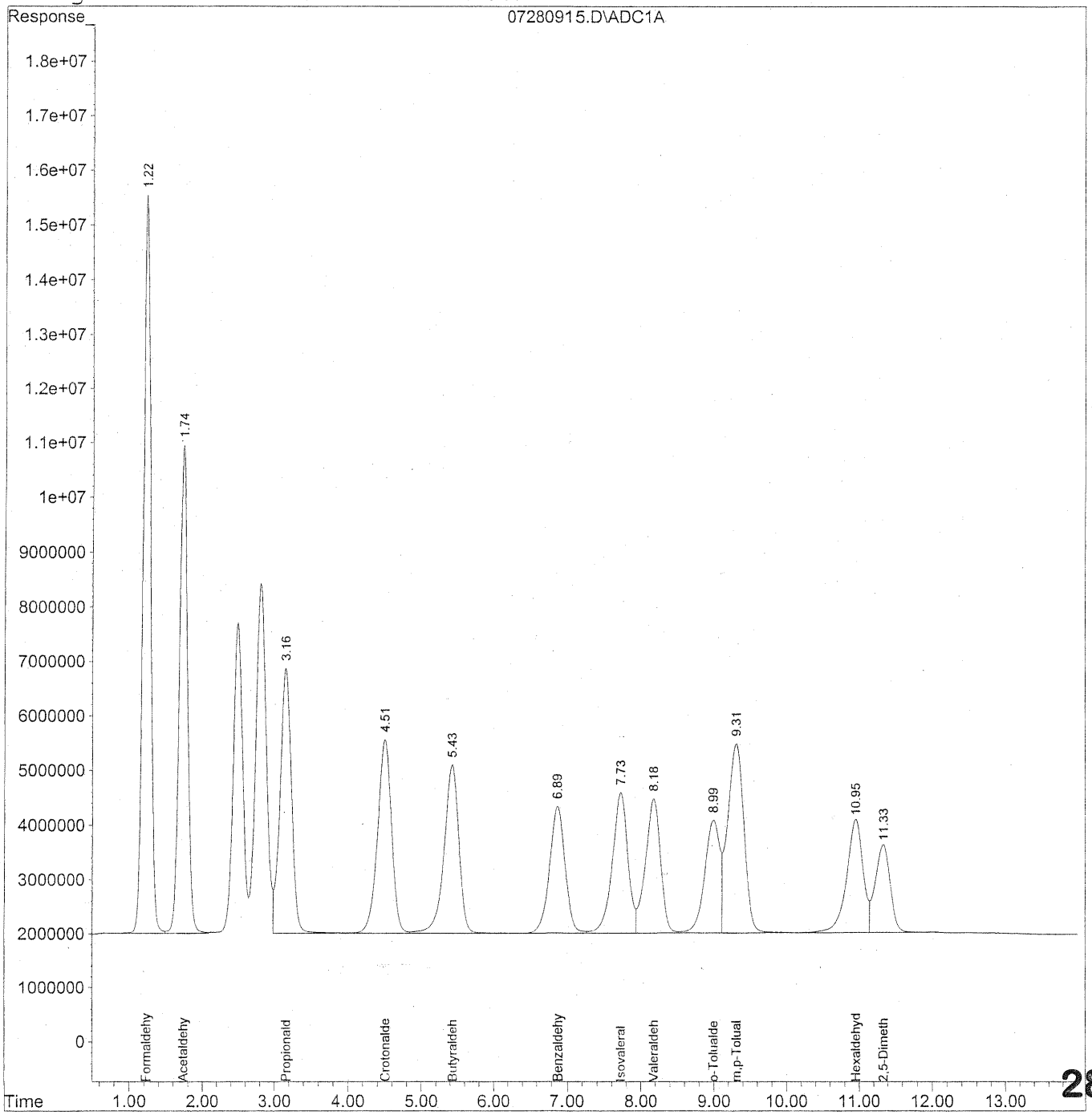
(f)=RT Delta > 1/2 Window (m)=manual int.

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



280

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

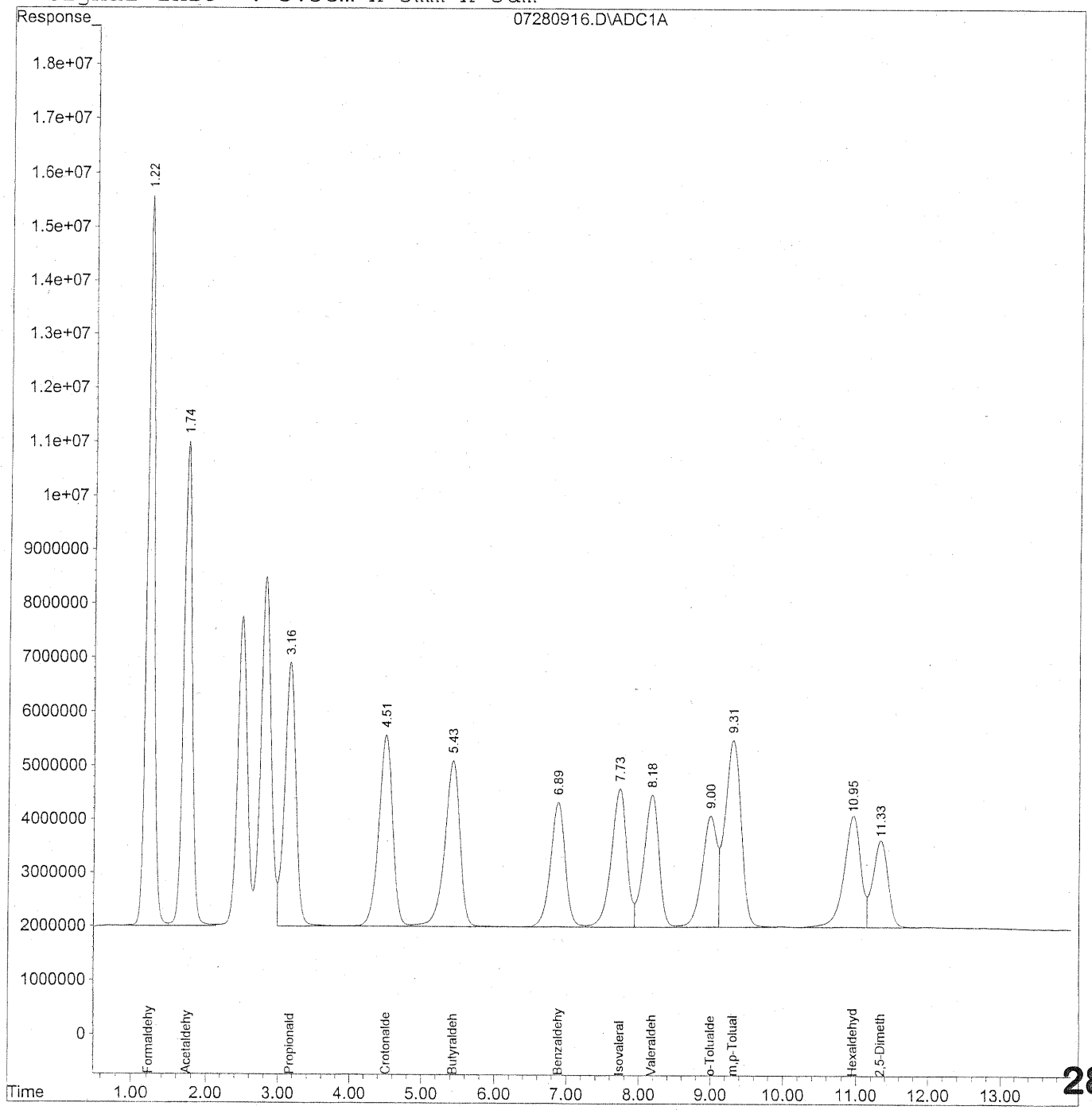
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	928364658	5175.655	ng/ml
2)	Acetaldehyde	1.74	706170560	5146.384	ng/ml
3)	Propionaldehyde	3.16	539067854	5164.924	ng/ml
4)	Crotonaldehyde	4.51	476268543	4639.802	ng/ml
5)	Butyraldehyde	5.43	446392739	5287.707	ng/ml
6)	Benzaldehyde	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.

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



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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

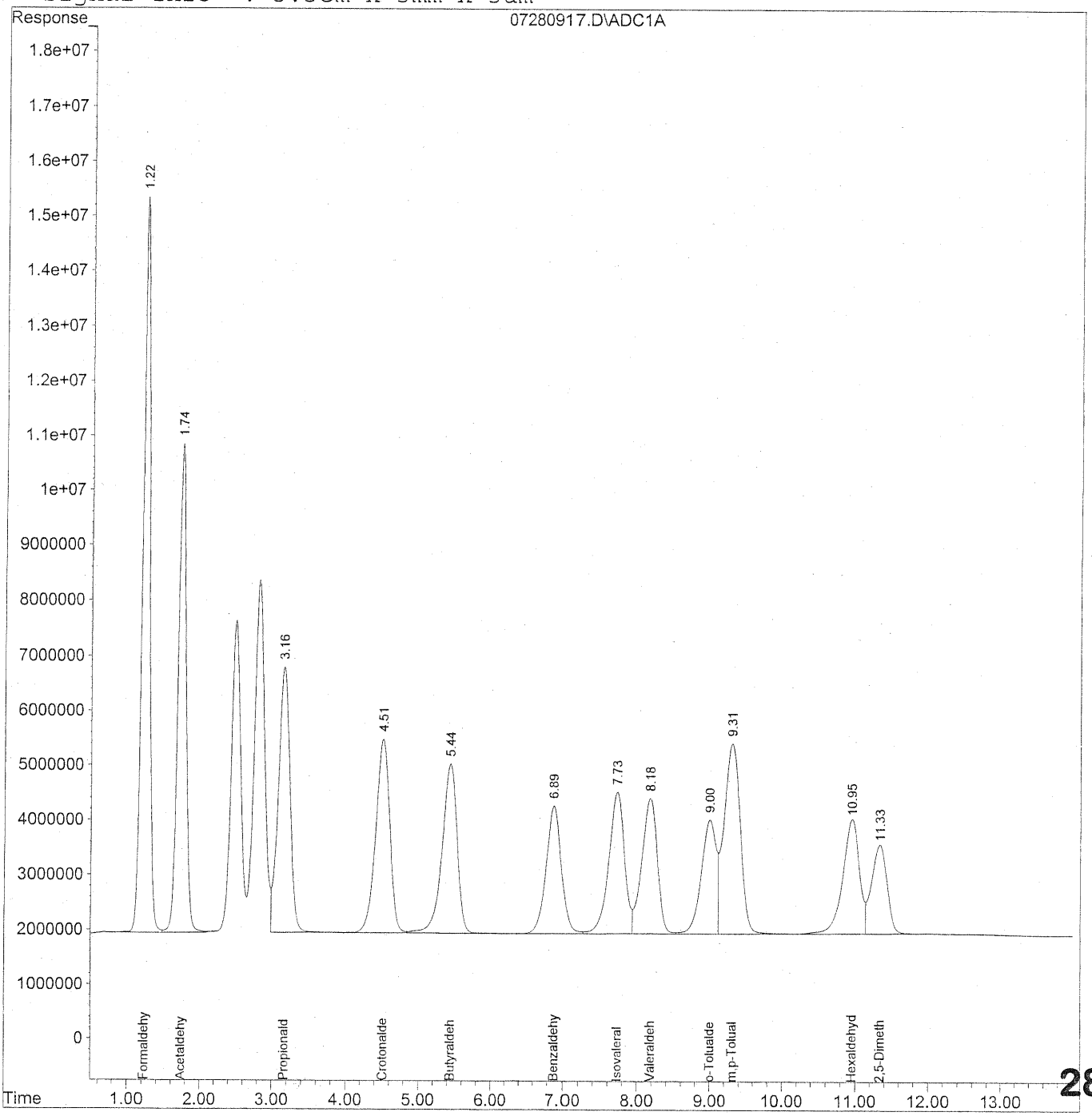
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	925768000	5161.179	ng/ml
2)	Acetaldehyde	1.74	708552415	5163.742	ng/ml
3)	Propionaldehyde	3.16	540133923	5175.139	ng/ml
4)	Crotonaldehyde	4.51	477844499	4655.155	ng/ml
5)	Butyraldehyde	5.43	446568052	5289.783	ng/ml
6)	Benzaldehyde	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.

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



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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

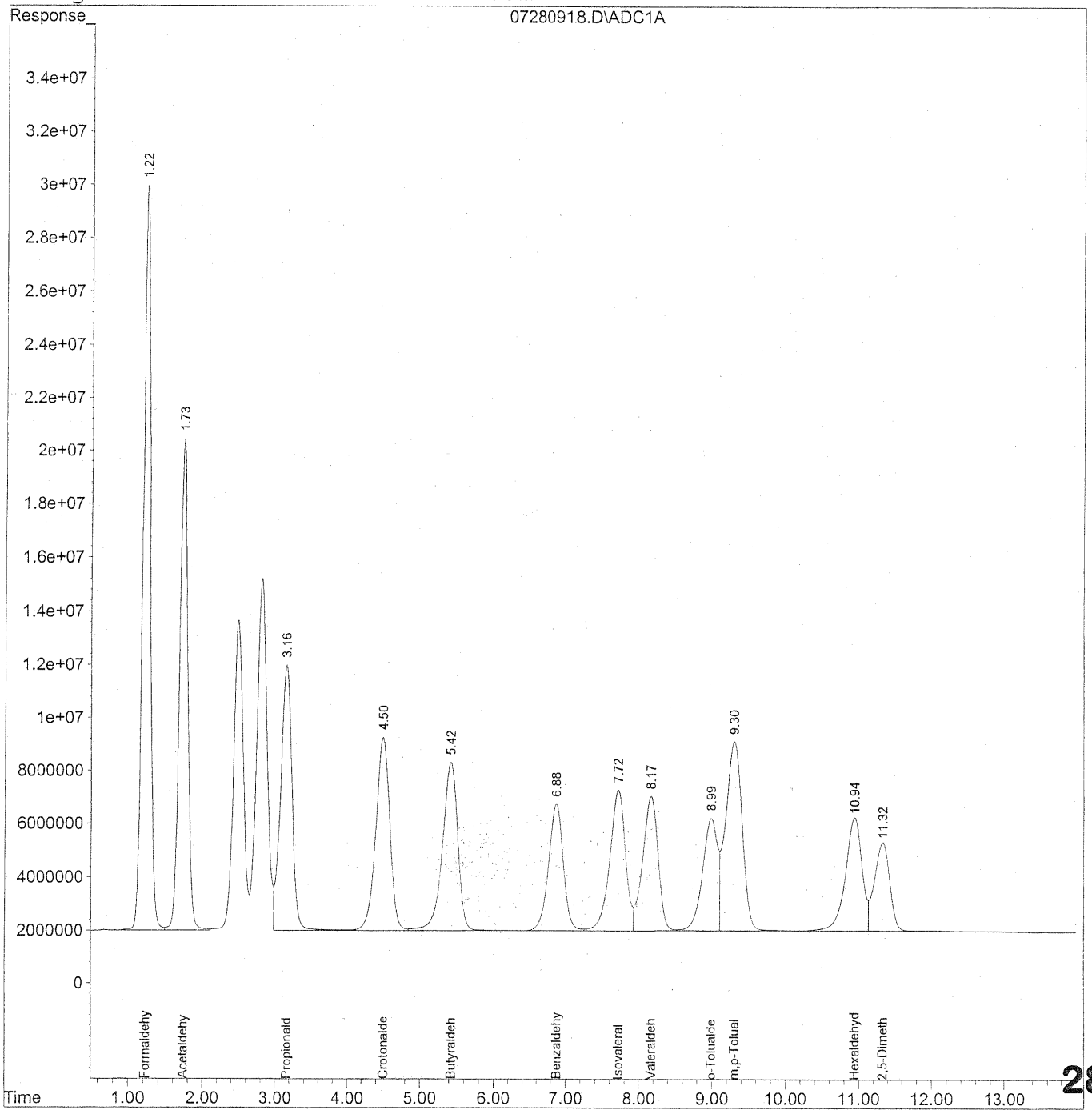
(f)=RT Delta > 1/2 Window (m)=manual int.

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



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

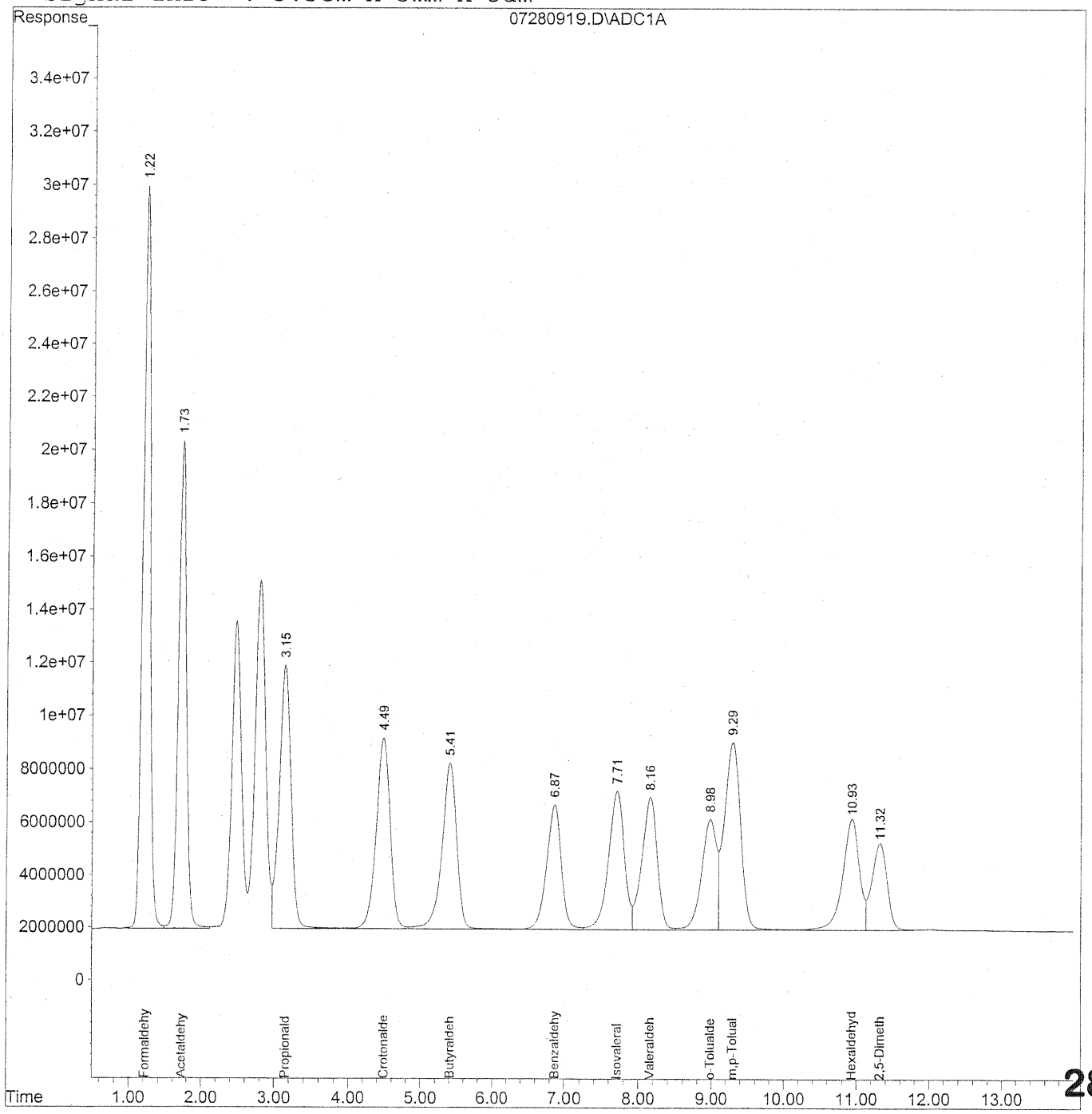
	Compound	R.T.	Response	Conc Units
	Target Compounds			
1)	Formaldehyde	1.22	1908653125	10640.787 ng/ml
2)	Acetaldehyde	1.73	1450154617	10568.343 ng/ml
3)	Propionaldehyde	3.16	1099941045	10538.770 ng/ml
4)	Crotonaldehyde	4.50	972691462	9475.947 ng/ml
5)	Butyraldehyde	5.42	910896701	10789.948 ng/ml
6)	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.

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



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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

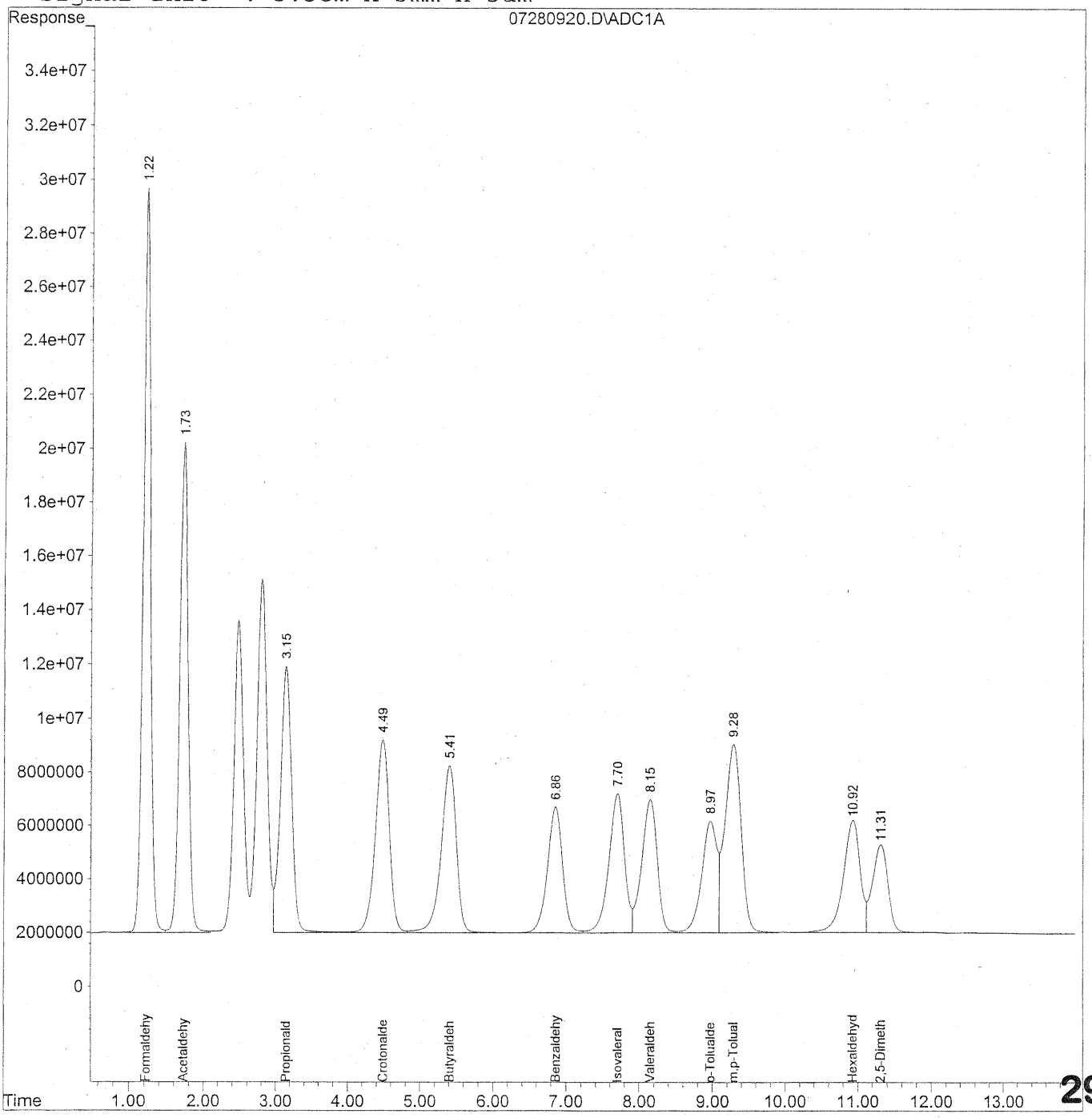
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	1905913073	10625.511	ng/ml
2)	Acetaldehyde	1.73	1446499891	10541.708	ng/ml
3)	Propionaldehyde	3.15	1098837646	10528.198	ng/ml
4)	Crotonaldehyde	4.49	971357788	9462.954	ng/ml
5)	Butyraldehyde	5.41	911328243	10795.060	ng/ml
6)	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.

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

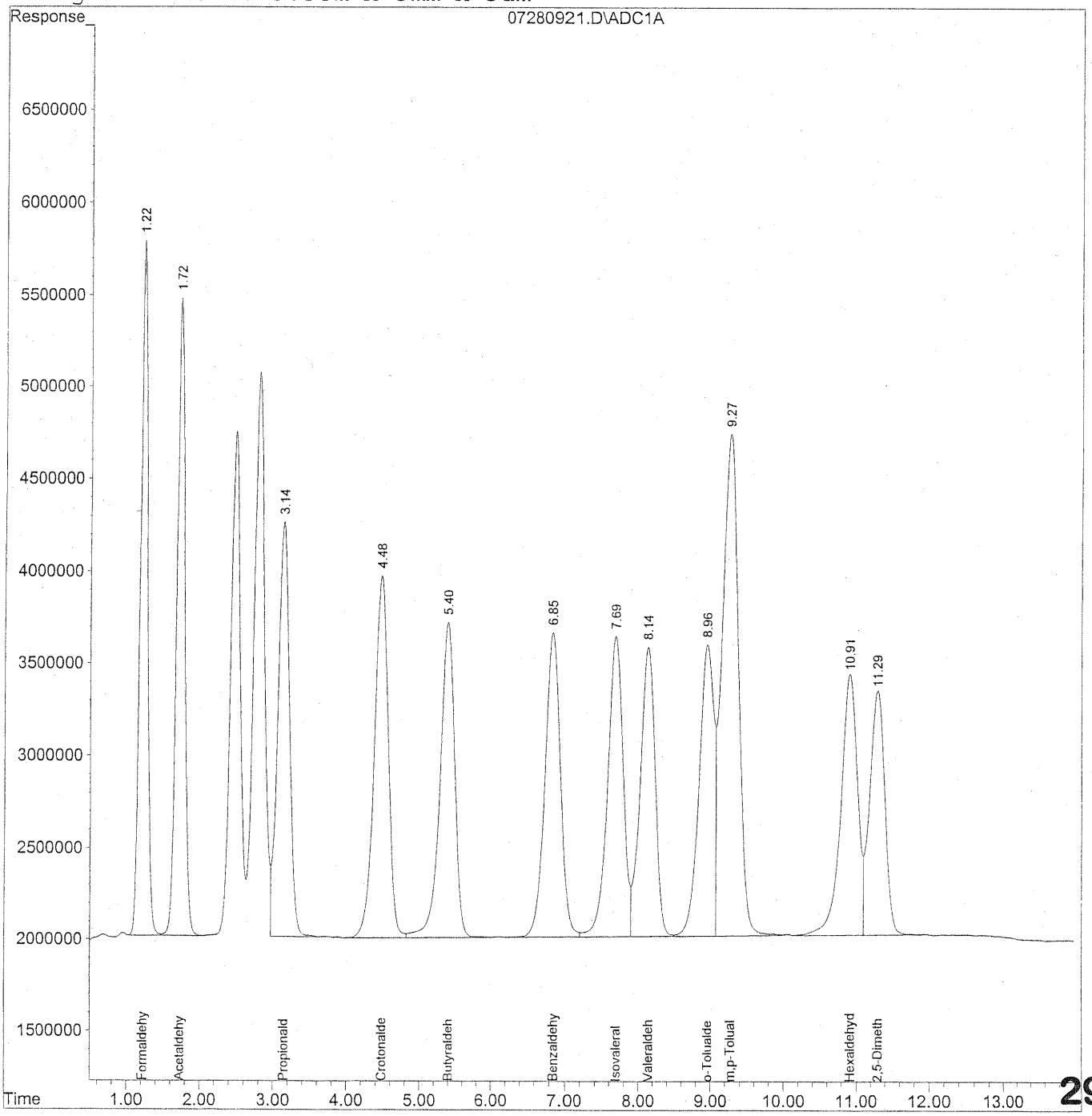
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	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

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



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

Printed : 8/31/09 hll 8/31/09

Instrument : LC#1 | Date Acquired : 8/26/09

Detector : UV-VIS 360 | Sample Amount : 5ul

Analyst : HC | Client & PAI Job# : EH&E P0902942

SAMPLE RESULT SUMMARY

Sample Information	MDL	1500ng/ml TO11A std S21-08250901	% Diff	ACN blank Lot CY023	MB front lot 5855/5994 1.0ml	MB back lot 5855/5994 1.0ml	P0902942-001 back 1.0ml	P0902942-002 back 1.0ml	P0902942-003 back 1.0ml
Dilution	1.0			1.0	1.0	1.0	1.0	1.0	1.0
Sample Volume (L)	NA			NA	NA	NA	104.04	104.04	100.50
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	ng/sample
Formaldehyde	100.00	1362.2	9.2%	ND	ND	ND	ND	ND	198.683 B,H
Acetaldehyde	100.00	1360.3	9.3%	ND	ND	ND	230.544	296.120	ND
Propionaldehyde	100.00	1347.6	10.2%	ND	ND	ND	ND	ND	ND
Crotonaldehyde	100.00	1323.8	11.7%	ND	ND	ND	ND	ND	ND
Butyraldehyde	100.00	1360.8	9.3%	ND	ND	ND	ND	ND	ND
Benzaldehyde	100.00	1328.5	11.4%	ND	ND	ND	ND	ND	ND
Isovaleraldehyde	100.00	1316.2	12.3%	ND	ND	ND	ND	ND	ND
Valeraldehyde	100.00	1300.6	13.3%	ND	ND	ND	ND	ND	ND
o-Tolualdehyde	100.00	1345.4	10.3%	ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde	200.00	2625.8	12.5%	ND	ND	ND	ND	ND	ND
Hexaldehyde	100.00	1366.2	8.9%	ND	ND	ND	ND	ND	ND
2,5-Dimethylbenzaldehyde	100.00	1281.0	14.6%	ND	ND	ND	ND	ND	ND

	ug/m3		ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3
Formaldehyde			NA	NA	NA	ND	ND	1.977
Acetaldehyde			NA	NA	NA	2.216	2.846	ND
Propionaldehyde			NA	NA	NA	ND	ND	ND
Crotonaldehyde			NA	NA	NA	ND	ND	ND
Butyraldehyde			NA	NA	NA	ND	ND	ND
Benzaldehyde			NA	NA	NA	ND	ND	ND
Isovaleraldehyde			NA	NA	NA	ND	ND	ND
Valeraldehyde			NA	NA	NA	ND	ND	ND
o-Tolualdehyde			NA	NA	NA	ND	ND	ND
m,p-Tolualdehyde			NA	NA	NA	ND	ND	ND
Hexaldehyde			NA	NA	NA	ND	ND	ND
2,5-Dimethylbenzaldehyde			NA	NA	NA	ND	ND	ND

	ppb		ppb	ppb	ppb	ppb	ppb	ppb
Formaldehyde			NA	NA	NA	ND	ND	1.610
Acetaldehyde			NA	NA	NA	1.230	1.580	ND
Propionaldehyde			NA	NA	NA	ND	ND	ND
Crotonaldehyde			NA	NA	NA	ND	ND	ND
Butyraldehyde			NA	NA	NA	ND	ND	ND
Benzaldehyde			NA	NA	NA	ND	ND	ND
Isovaleraldehyde			NA	NA	NA	ND	ND	ND
Valeraldehyde			NA	NA	NA	ND	ND	ND
o-Tolualdehyde			NA	NA	NA	ND	ND	ND
m,p-Tolualdehyde			NA	NA	NA	ND	ND	ND
Hexaldehyde			NA	NA	NA	ND	ND	ND
2,5-Dimethylbenzaldehyde			NA	NA	NA	ND	ND	ND

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

Printed : 8/31/09
Instrument : LC#1 Date Acquired : 8/26/09
Detector : UV-VIS 360 Sample Amount : 5ul
Analyst : HC Client & PAI Job# : EH&E P0902942

SAMPLE RESULT SUMMARY

Sample Information	MDL	P0902942-004 back 1.0ml	P0902942-005 back 1.0ml	P0902942-006 back 1.0ml	CCV 1500ng/ml S21-08250901	% Diff	CCV 1500ng/ml S21-08250901	% Diff
Dilution	1.0	1.0	1.0	1.0	1.0		1.0	
Sample Volume (L)	NA	101.50	104.54	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	
Formaldehyde	100.00	ND	ND	ND	1434.109	4.4%	1398.642	6.8%
Acetaldehyde	100.00	270.341	259.749	ND	1413.822	5.7%	1398.945	6.7%
Propionaldehyde	100.00	ND	ND	ND	1410.148	6.0%	1383.422	7.8%
Crotonaldehyde	100.00	ND	ND	ND	1377.245	8.2%	1353.012	9.8%
Butyraldehyde	100.00	ND	ND	ND	1429.192	4.7%	1420.016	5.3%
Benzaldehyde	100.00	ND	ND	ND	1437.436	4.2%	1403.549	6.4%
Isovaleraldehyde	100.00	ND	ND	ND	1451.667	3.2%	1430.935	4.6%
Valeraldehyde	100.00	ND	ND	ND	1382.046	7.9%	1328.692	11.4%
o-Tolualdehyde	100.00	ND	ND	ND	1481.082	1.3%	1467.914	2.1%
m,p-Tolualdehyde	200.00	ND	ND	ND	2878.668	4.0%	2826.510	5.8%
Hexaldehyde	100.00	ND	ND	ND	1444.556	3.7%	1416.960	5.5%
2,5-Dimethylbenzaldehyde	100.00	ND	ND	ND	1286.565	14.2%	1300.850	13.3%

	ug/m3	ug/m3	ug/m3
Formaldehyde	ND	ND	ND
Acetaldehyde	2.663	2.485	ND
Propionaldehyde	ND	ND	ND
Crotonaldehyde	ND	ND	ND
Butyraldehyde	ND	ND	ND
Benzaldehyde	ND	ND	ND
Isovaleraldehyde	ND	ND	ND
Valeraldehyde	ND	ND	ND
o-Tolualdehyde	ND	ND	ND
m,p-Tolualdehyde	ND	ND	ND
Hexaldehyde	ND	ND	ND
2,5-Dimethylbenzaldehyde	ND	ND	ND

	ppb	ppb	ppb
Formaldehyde	ND	ND	ND
Acetaldehyde	1.479	1.380	ND
Propionaldehyde	ND	ND	ND
Crotonaldehyde	ND	ND	ND
Butyraldehyde	ND	ND	ND
Benzaldehyde	ND	ND	ND
Isovaleraldehyde	ND	ND	ND
Valeraldehyde	ND	ND	ND
o-Tolualdehyde	ND	ND	ND
m,p-Tolualdehyde	ND	ND	ND
Hexaldehyde	ND	ND	ND
2,5-Dimethylbenzaldehyde	ND	ND	ND

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

Instrument : LC#1
Detector : UV-VIS 360
Analyst : HC

Printed : 8/31/09
Date Acquired : 8/26/09
Sample Amount : 5ul
Client & PAI Job# : EH&E P0902942

Sample Information	MDL	ACN lot blk CY023	MB front lot 5855/5994 1.0ml	MB back lot 5855/5994 1.0ml	CCV 1500ng/ml S21-08250901	% Diff	P0902942-001 front 1.0ml	P0902942-001 front 1.0ml
Dilution	1.0	1.0	1.0	1.0	1.0		1.0	1.0
Sample Volume (L)	NA						104.04	104.04
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	1398.621	6.8%	2319.043	2319.043
Acetaldehyde	100.00	ND	ND	ND	1401.357	6.6%	4377.863	4377.863
Propionaldehyde	100.00	ND	ND	ND	1380.821	7.9%	594.458	594.458
Crotonaldehyde	100.00	ND	ND	ND	1358.913	9.4%	ND	ND
Butyraldehyde	100.00	ND	ND	ND	1397.881	6.8%	1046.320	1046.320
Benzaldehyde	100.00	ND	ND	ND	1381.176	7.9%	956.648	956.648
Isovaleraldehyde	100.00	ND	ND	ND	1396.480	6.9%	154.901	154.901
Valeraldehyde	100.00	ND	ND	ND	1329.746	11.4%	1614.993	1614.993
o-Tolualdehyde	100.00	ND	ND	ND	1409.945	6.0%	ND	ND
m,p-Tolualdehyde	200.00	ND	ND	ND	2785.028	7.2%	ND	ND
Hexaldehyde	100.00	ND	ND	ND	1416.258	5.6%	4523.821	4523.821
2,5-Dimethylbenzaldehyde	100.00	ND	ND	ND	1296.689	13.6%	ND	ND

	ug/m3	ug/m3	ug/m3			ug/m3	ug/m3
Formaldehyde		ND	ND	ND		22.290	22.290
Acetaldehyde		ND	ND	ND		42.079	42.079
Propionaldehyde		ND	ND	ND		5.714	5.714
Crotonaldehyde		ND	ND	ND		ND	ND
Butyraldehyde		ND	ND	ND		10.057	10.057
Benzaldehyde		ND	ND	ND		9.195	9.195
Isovaleraldehyde		ND	ND	ND		1.489	1.489
Valeraldehyde		ND	ND	ND		15.523	15.523
o-Tolualdehyde		ND	ND	ND		ND	ND
m,p-Tolualdehyde		ND	ND	ND		ND	ND
Hexaldehyde		ND	ND	ND		43.482	43.482
2,5-Dimethylbenzaldehyde		ND	ND	ND		ND	ND

	ppb	ppb	ppb	ppb		ppb	ppb
Formaldehyde		ND	ND	ND		18.156	18.156
Acetaldehyde		ND	ND	ND		23.365	23.365
Propionaldehyde		ND	ND	ND		2.406	2.406
Crotonaldehyde		ND	ND	ND		ND	ND
Butyraldehyde		ND	ND	ND		3.411	3.411
Benzaldehyde		ND	ND	ND		2.119	2.119
Isovaleraldehyde		ND	ND	ND		0.423	0.423
Valeraldehyde		ND	ND	ND		4.408	4.408
o-Tolualdehyde		ND	ND	ND		ND	ND
m,p-Tolualdehyde		ND	ND	ND		ND	ND
Hexaldehyde		ND	ND	ND		10.619	10.619
2,5-Dimethylbenzaldehyde		ND	ND	ND		ND	ND

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

Instrument : LC#1
Detector : UV-VIS 360
Analyst : HC

Printed : 8/31/09
Date Acquired : 8/26/09
Sample Amount 5ul
Client & PAI Job EH&E P0902942

Sample Information	MDL	CCV 1500ng/ml S21-08250901	% Diff	1500ng/ml TO11A std S21-08270903	% Diff	ACN blank Lot CY023	MB front lot 5855/5994 1.0ml	MB back lot 5855/5994 1.0ml
Dilution	1.0	1.0		1.0		1.0	1.0	1.0
Sample Volume (L)	NA							
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
Formaldehyde	100.00	1408.687	6.1%	1313.306	12.4%	ND	ND	ND
Acetaldehyde	100.00	1384.061	7.7%	1304.154	13.1%	ND	ND	ND
Propionaldehyde	100.00	1370.631	8.6%	1290.921	13.9%	ND	ND	ND
Crotonaldehyde	100.00	1365.704	9.0%	1276.630	14.9%	ND	ND	ND
Butyraldehyde	100.00	1395.582	7.0%	1318.473	12.1%	ND	ND	ND
Benzaldehyde	100.00	1392.170	7.2%	1302.108	13.2%	ND	ND	ND
Isovaleraldehyde	100.00	1409.573	6.0%	1342.282	10.5%	ND	ND	ND
Valeraldehyde	100.00	1305.773	12.9%	1277.949	14.8%	ND	ND	ND
o-Tolualdehyde	100.00	1412.688	5.8%	1361.756	9.2%	ND	ND	ND
m,p-Tolualdehyde	200.00	2758.587	8.0%	2588.156	13.7%	ND	ND	ND
Hexaldehyde	100.00	1410.269	6.0%	1350.526	10.0%	ND	ND	ND
2,5-Dimethylbenzaldehyde	100.00	1297.167	13.5%	1232.334	17.8%	ND	ND	ND
	ug/m3					ug/m3	ug/m3	ug/m3
Formaldehyde						ND	ND	ND
Acetaldehyde						ND	ND	ND
Propionaldehyde						ND	ND	ND
Crotonaldehyde						ND	ND	ND
Butyraldehyde						ND	ND	ND
Benzaldehyde						ND	ND	ND
Isovaleraldehyde						ND	ND	ND
Valeraldehyde						ND	ND	ND
o-Tolualdehyde						ND	ND	ND
m,p-Tolualdehyde						ND	ND	ND
Hexaldehyde						ND	ND	ND
2,5-Dimethylbenzaldehyde						ND	ND	ND
	ppb					ppb	ppb	ppb
Formaldehyde						ND	ND	ND
Acetaldehyde						ND	ND	ND
Propionaldehyde						ND	ND	ND
Crotonaldehyde						ND	ND	ND
Butyraldehyde						ND	ND	ND
Benzaldehyde						ND	ND	ND
Isovaleraldehyde						ND	ND	ND
Valeraldehyde						ND	ND	ND
o-Tolualdehyde						ND	ND	ND
m,p-Tolualdehyde						ND	ND	ND
Hexaldehyde						ND	ND	ND
2,5-Dimethylbenzaldehyde						ND	ND	ND

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

Instrument LC#1
Detector : UV-VIS 360
Analyst : HC

Printed : 8/31/09
Date Acquirec 8/26/09
Sample Amou 5ul
Client & PAI J EH&E P0902942

Sample Information	MDL	P0902942-002 front 1.0ml	P0902942-003 front 1.0ml	P0902942-004 front 1.0ml	P0902942-005 front 1.0ml	P0902942-006 front 1.0ml	CCV 1500ng/ml S21-08270903	% Diff
Dilution	1.0	1.0	1.0	1.0	1.0	1.0	1.0	
Sample Volume (L)	NA	104.04	100.50	101.50	104.54	0.0		
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	2325.872	ND	2119.457	2268.907	ND	1289.375	14.0%
Acetaldehyde	100.00	4444.588	113.231	4072.128	4532.977	ND	1281.442	14.6%
Propionaldehyde	100.00	567.173	ND	539.869	635.193	ND	1282.582	14.5%
Crotonaldehyde	100.00	ND	ND	ND	ND	ND	1255.434	16.3% V
Butyraldehyde	100.00	1070.970 MP	ND	1007.640	1100.646 MP	ND	1282.940	14.5%
Benzaldehyde	100.00	839.526	ND	820.072	796.082	ND	1283.728	14.4%
Isovaleraldehyde	100.00	171.663	ND	167.665	137.045	ND	1340.786	10.6%
Valeraldehyde	100.00	1536.409	ND	1440.994	1430.644	ND	1277.726	14.8%
o-Tolualdehyde	100.00	ND	ND	ND	ND	ND	1353.149	9.8%
m,p-Tolualdehyde	200.00	ND	ND	ND	ND	ND	2636.438	12.1%
Hexaldehyde	100.00	4578.303	ND	4193.832	5021.234	ND	1521.926	1.5%
2,5-Dimethylbenzaldehyde	100.00	ND	ND	ND	ND	ND	1518.971	1.3%

	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3
Formaldehyde		22.356	ND	20.881	21.704	ND
Acetaldehyde		42.720	1.127	40.119	43.361	ND
Propionaldehyde		5.451	ND	5.319	6.076	ND
Crotonaldehyde		ND	ND	ND	ND	ND
Butyraldehyde		10.294	ND	9.927	10.528	ND
Benzaldehyde		8.069	ND	8.080	7.615	ND
Isovaleraldehyde		1.650	ND	1.652	1.311	ND
Valeraldehyde		14.767	ND	14.197	13.685	ND
o-Tolualdehyde		ND	ND	ND	ND	ND
m,p-Tolualdehyde		ND	ND	ND	ND	ND
Hexaldehyde		44.005	ND	41.319	48.032	ND
2,5-Dimethylbenzaldehyde		ND	ND	ND	ND	ND

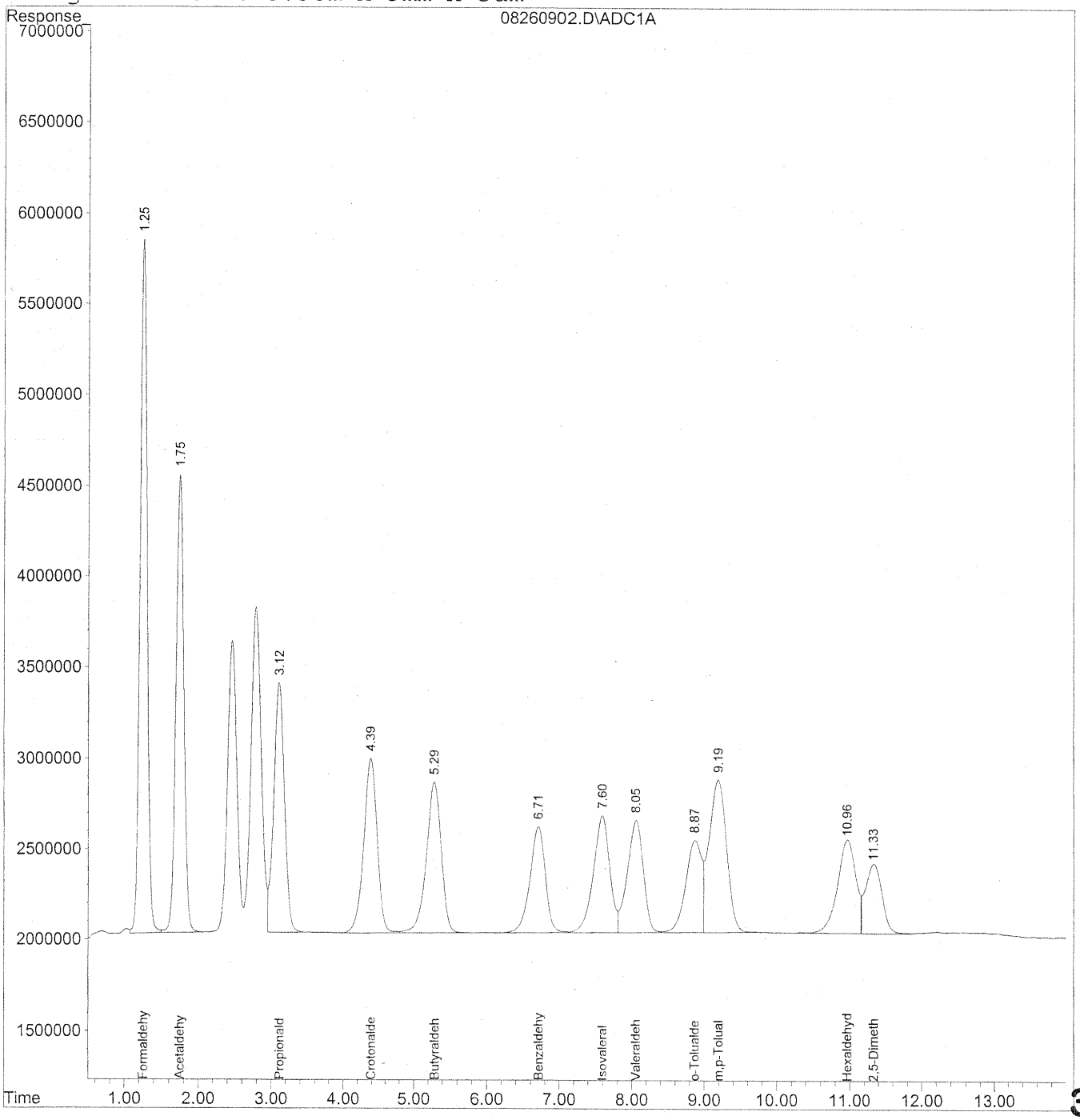
	ppb	ppb	ppb	ppb	ppb	ppb
Formaldehyde		18.209	ND	17.008	17.678	ND
Acetaldehyde		23.721	0.626	22.277	24.078	ND
Propionaldehyde		2.296	ND	2.240	2.559	ND
Crotonaldehyde		ND	ND	ND	ND	ND
Butyraldehyde		3.492	ND	3.367	3.571	ND
Benzaldehyde		1.860	ND	1.862	1.755	ND
Isovaleraldehyde		0.469	ND	0.469	0.372	ND
Valeraldehyde		4.194	ND	4.032	3.886	ND
o-Tolualdehyde		ND	ND	ND	ND	ND
m,p-Tolualdehyde		ND	ND	ND	ND	ND
Hexaldehyde		10.746	ND	10.090	11.730	ND
2,5-Dimethylbenzaldehyde		ND	ND	ND	ND	ND

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260902.D Vial: 2
Acq On : 26 Aug 2009 5:20 pm Operator: HC
Sample : 1500ng/ml TO11A std S21-08250901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:30 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 : Sat Aug 29 12:41:27 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\26\08260902.D Vial: 2
Acq On : 26 Aug 2009 5:20 pm Operator: HC
Sample : 1500ng/ml TO11A std S21-08250901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:30 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 : Sat Aug 29 12:41:27 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

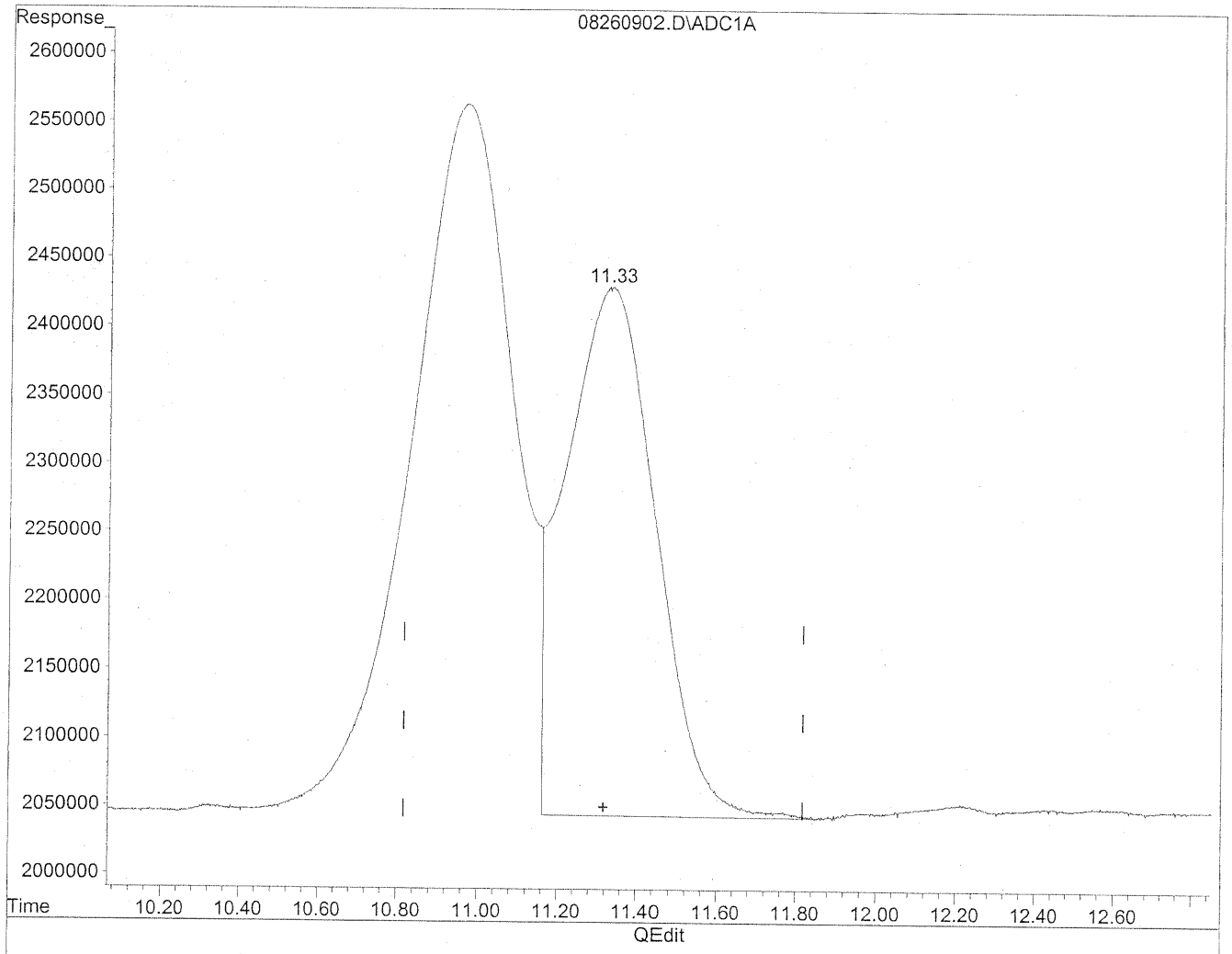
8/29/07

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.25	250073526	1362.194	ng/ml
2)	Acetaldehyde	1.75	190751677	1360.340	ng/ml
3)	Propionaldehyde	3.12	143783952	1347.614	ng/ml
4)	Crotonaldehyde	4.39	128959783	1323.816	ng/ml
5)	Butyraldehyde	5.29	120211032	1360.837	ng/ml
6)	Benzaldehyde	6.71	87506049	1328.480	ng/ml
7)	Isovaleraldehyde	7.59	102992522	1316.182	ng/ml
8)	Valeraldehyde	8.06	95599983	1300.591	ng/ml
9)	o-Tolualdehyde	8.87	78465289	1345.417	ng/ml
10)	m,p-Tolualdehyde	9.19	141781732	2625.811	ng/ml
11)	Hexaldehyde	10.97	92006495	1366.221	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.33	62784813	1280.973	ng/mlm

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260902.D Vial: 2
Acq On : 26 Aug 2009 5:20 pm Operator: HC
Sample : 1500ng/ml TO11A std S21-08250901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:29 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 : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

11.33min 1246.963ng/ml

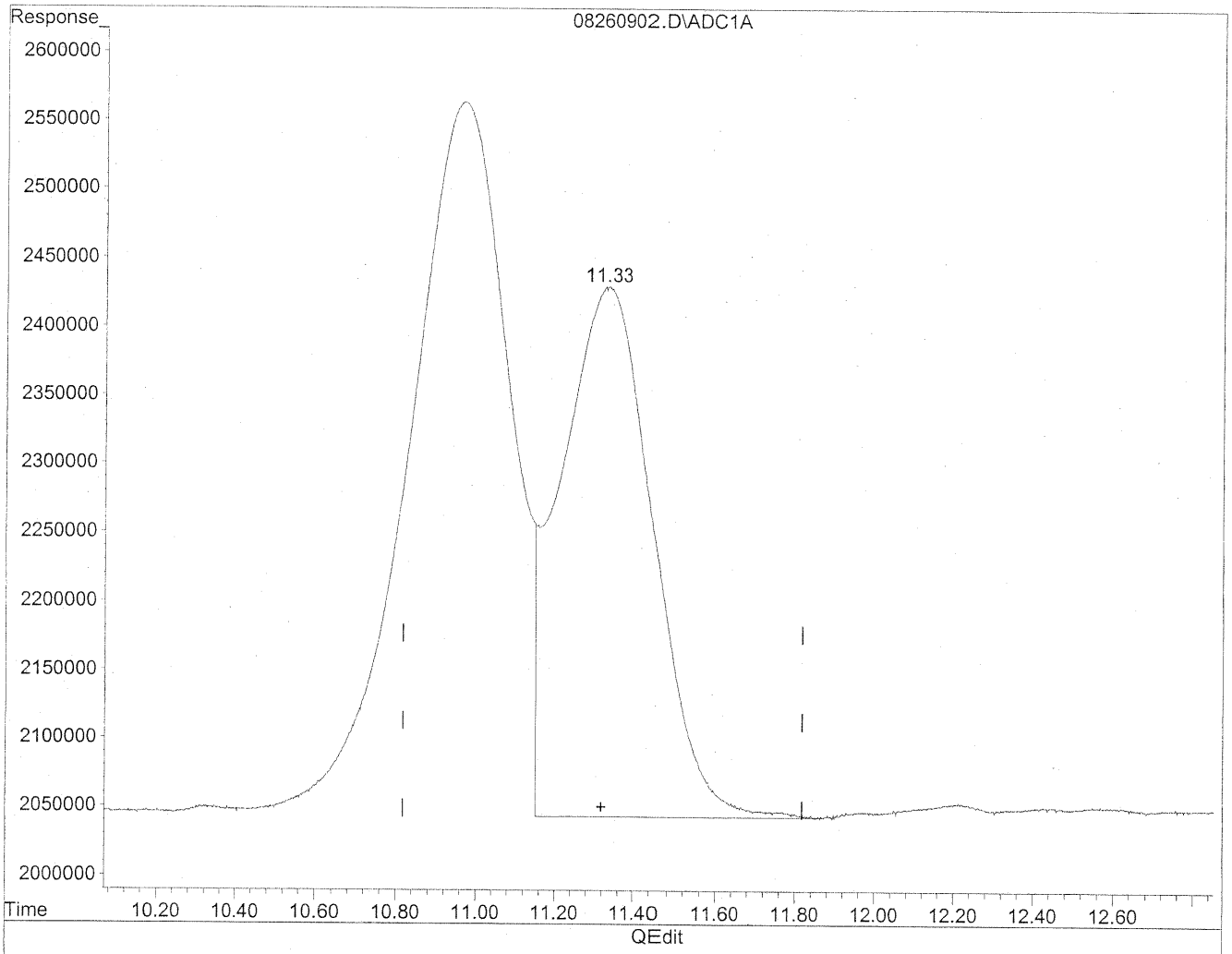
response 61117876

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260902.D Vial: 2
Acq On : 26 Aug 2009 5:20 pm Operator: HC
Sample : 1500ng/ml TO11A std S21-08250901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:29 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 : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

11.33min 1280.973ng/ml m

response 62784813

(+) = Expected Retention Time

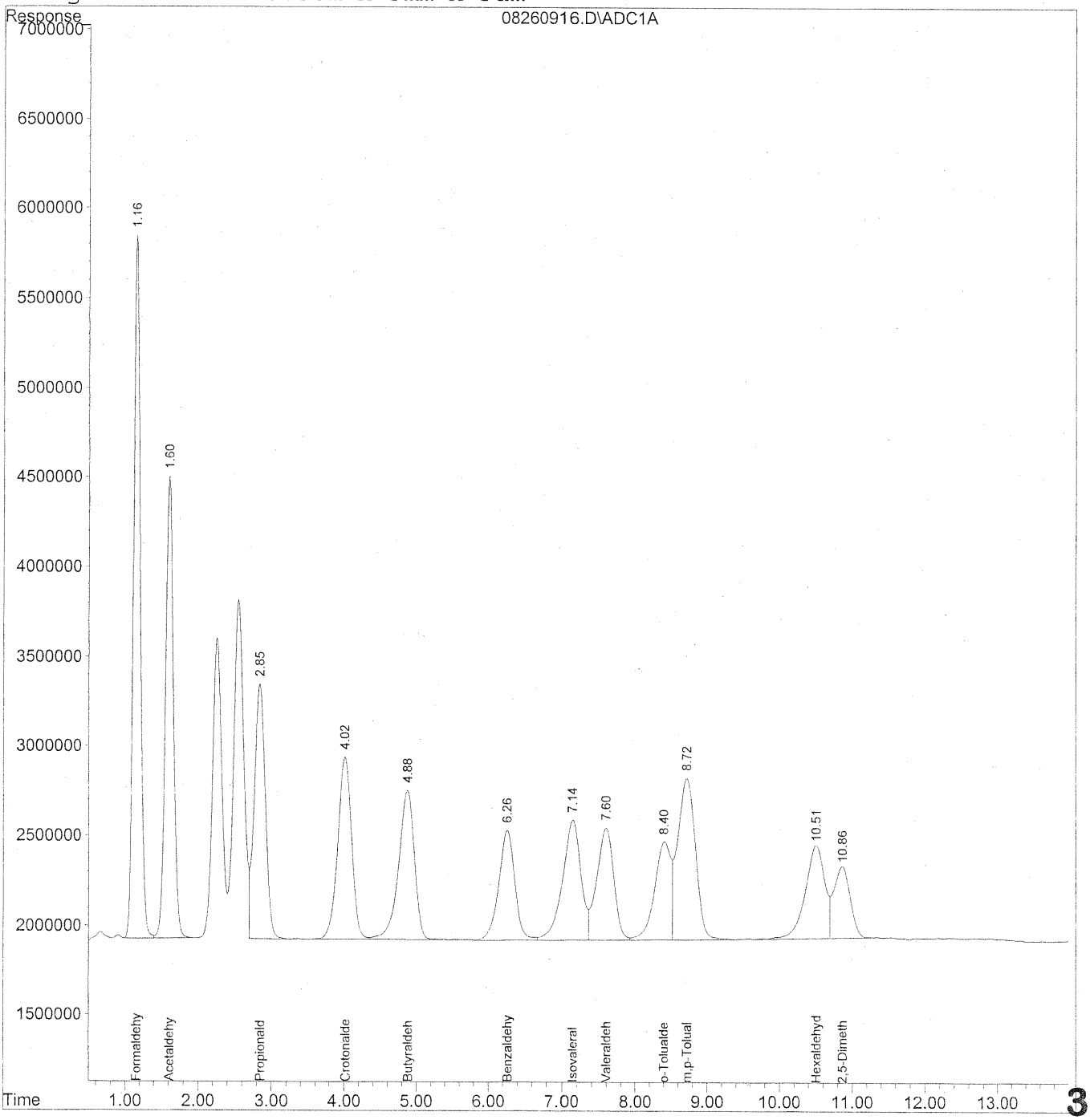
08260902.D TO110709.M Sat Aug 29 16:30:32 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260916.D Vial: 16
Acq On : 26 Aug 2009 8:51 pm Operator: HC
Sample : CCV 1500ng/ml S21-08250901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:33 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 : Sat Aug 29 16:33:38 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\26\08260916.D Vial: 16
Acq On : 26 Aug 2009 8:51 pm Operator: HC
Sample : CCV 1500ng/ml S21-08250901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 16:33 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 : Sat Aug 29 16:33:38 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	263275750	1434.109	ng/ml
2)	Acetaldehyde	1.61	198251029	1413.822	ng/ml
3)	Propionaldehyde	2.85	150456040	1410.148	ng/ml
4)	Crotonaldehyde	4.02	134164517	1377.245	ng/ml
5)	Butyraldehyde	4.88	126249251	1429.192	ng/ml
6)	Benzaldehyde	6.27	94682891	1437.436	ng/ml
7)	Isovaleraldehyde	7.15	113594414	1451.667	ng/ml
8)	Valeraldehyde	7.60	101587273	1382.046	ng/ml
9)	o-Tolualdehyde	8.41	86377303	1481.082	ng/ml
10)	m,p-Tolualdehyde	8.72	155434812	2878.668	ng/ml
11)	Hexaldehyde	10.50	97281818	1444.556	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.86	63058911	1286.565	ng/ml

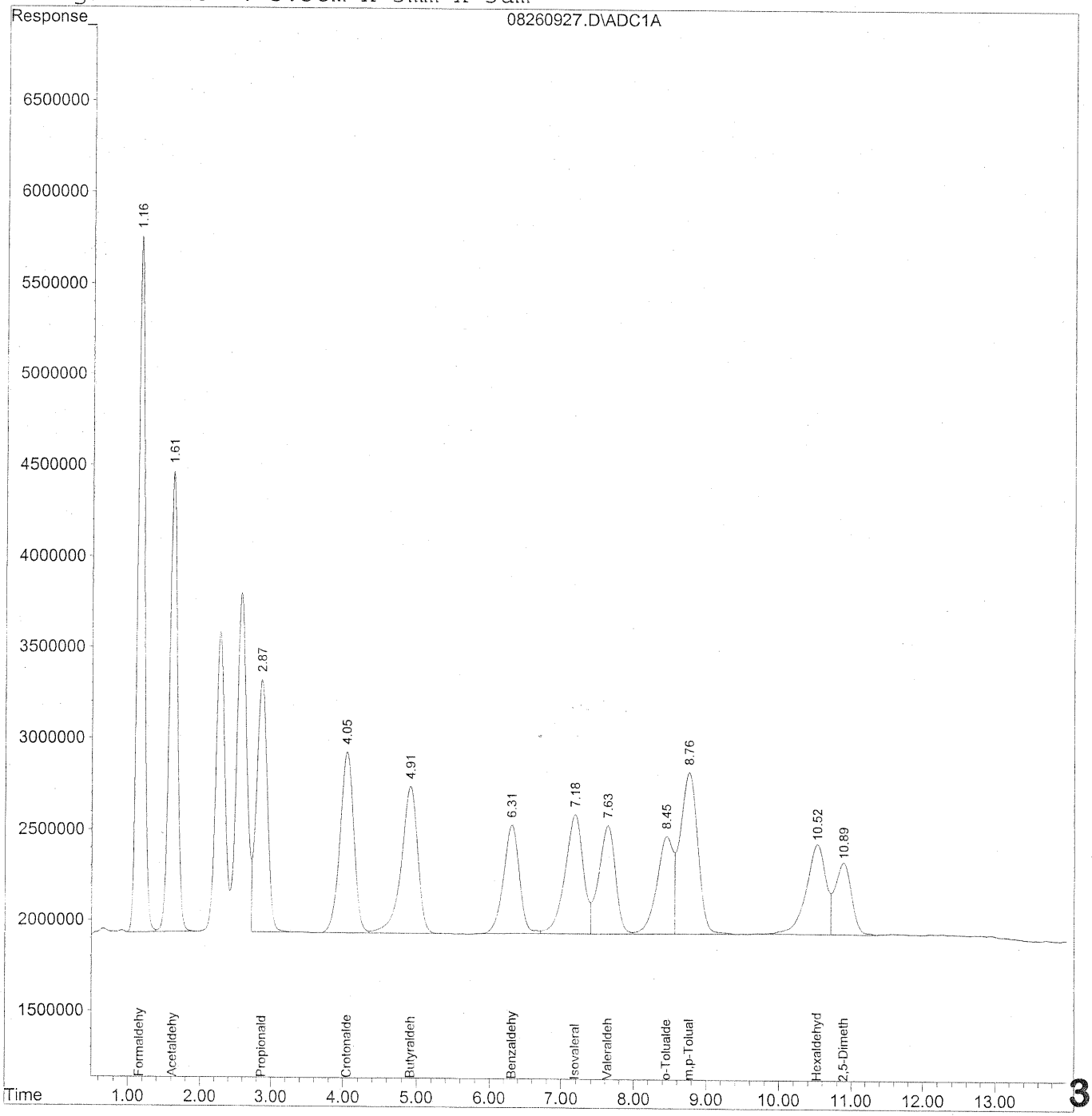
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260927.D Vial: 26
Acq On : 26 Aug 2009 11:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08250901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:36 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 : Thu Aug 27 07:35:56 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\26\08260927.D Vial: 26
Acq On : 26 Aug 2009 11:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08250901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:36 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 : Thu Aug 27 07:35:56 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

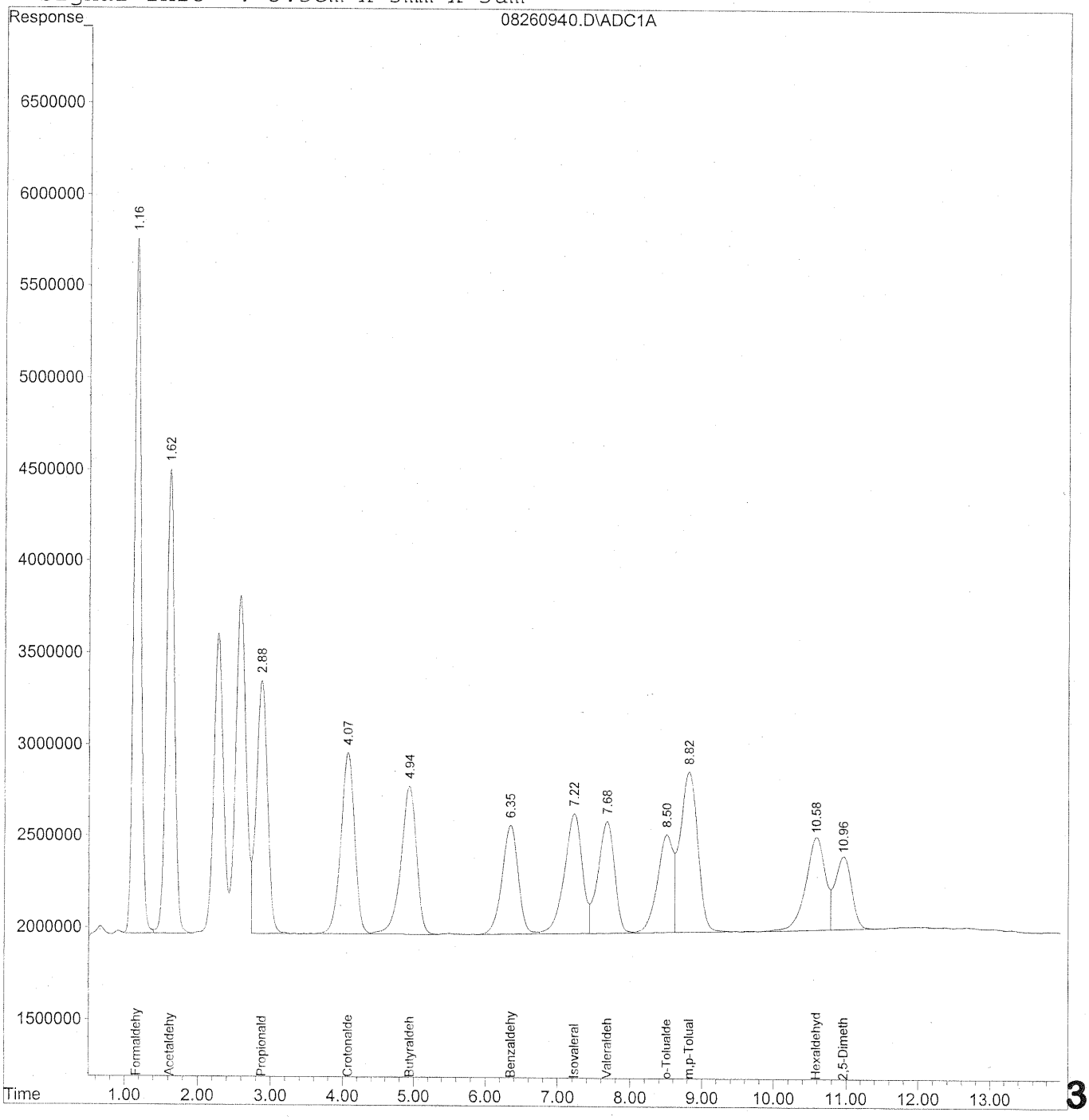
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	256764708	1398.642	ng/ml
2)	Acetaldehyde	1.61	196165014	1398.945	ng/ml
3)	Propionaldehyde	2.87	147604534	1383.422	ng/ml
4)	Crotonaldehyde	4.05	131803903	1353.012	ng/ml
5)	Butyraldehyde	4.91	125438742	1420.016	ng/ml
6)	Benzaldehyde	6.31	92450795	1403.549	ng/ml
7)	Isovaleraldehyde	7.18	111972071	1430.935	ng/ml
8)	Valeraldehyde	7.63	97665537	1328.692	ng/ml
9)	o-Tolualdehyde	8.45	85609330	1467.914	ng/ml
10)	m,p-Tolualdehyde	8.76	152618537	2826.510	ng/ml
11)	Hexaldehyde	10.52	95423426	1416.960	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.89	63759087	1300.850	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\26\08260940.D Vial: 39
Acq On : 27 Aug 2009 2:51 am Operator: HC
Sample : CCV 1500ng/ml S21-08250901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:36 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 : Thu Aug 27 07:35:56 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



309

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\26\08260940.D Vial: 39
Acq On : 27 Aug 2009 2:51 am Operator: HC
Sample : CCV 1500ng/ml S21-08250901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:36 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 : Thu Aug 27 07:35:56 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

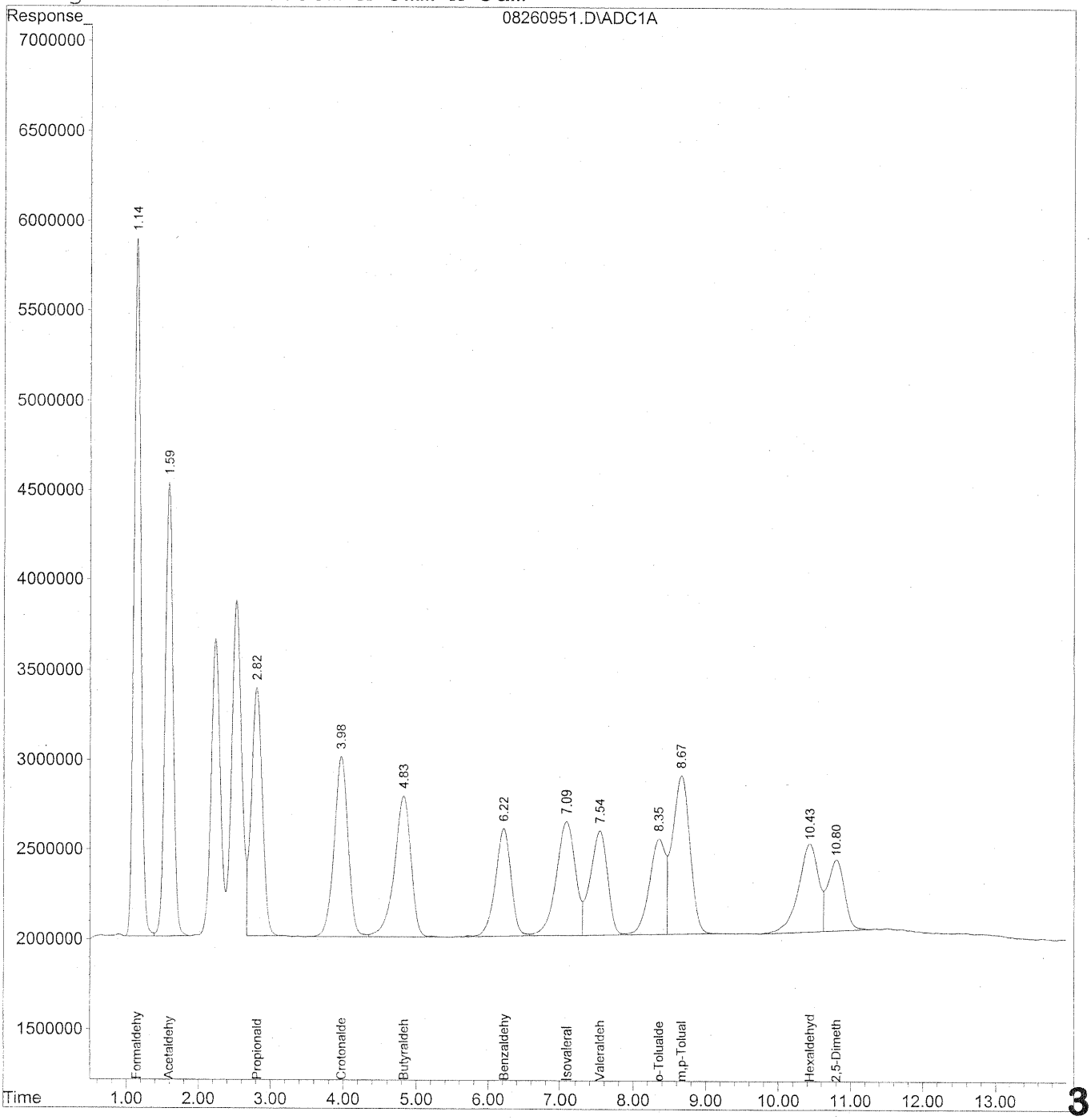
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.16	256760803	1398.621	ng/ml
2)	Acetaldehyde	1.62	196503156	1401.357	ng/ml
3)	Propionaldehyde	2.88	147327009	1380.821	ng/ml
4)	Crotonaldehyde	4.07	132378738	1358.913	ng/ml
5)	Butyraldehyde	4.94	123483374	1397.881	ng/ml
6)	Benzaldehyde	6.35	90977117	1381.176	ng/ml
7)	Isovaleraldehyde	7.22	109275943	1396.480	ng/ml
8)	Valeraldehyde	7.68	97742964	1329.746	ng/ml
9)	o-Tolualdehyde	8.50	82228602	1409.945	ng/ml
10)	m,p-Tolualdehyde	8.82	150378717	2785.028	ng/ml
11)	Hexaldehyde	10.59	95376129	1416.258	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.96	63555119	1296.689	ng/ml

Data File : J:\LC01\DATA\TO11\2009_08\26\08260951.D Vial: 49
Acq On : 27 Aug 2009 5:37 am Operator: HC
Sample : CCV 1500ng/ml S21-08250901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:36 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 : Thu Aug 27 07:35:56 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\26\08260951.D Vial: 49
Acq On : 27 Aug 2009 5:37 am Operator: HC
Sample : CCV 1500ng/ml S21-08250901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 7:36 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 : Thu Aug 27 07:35:56 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

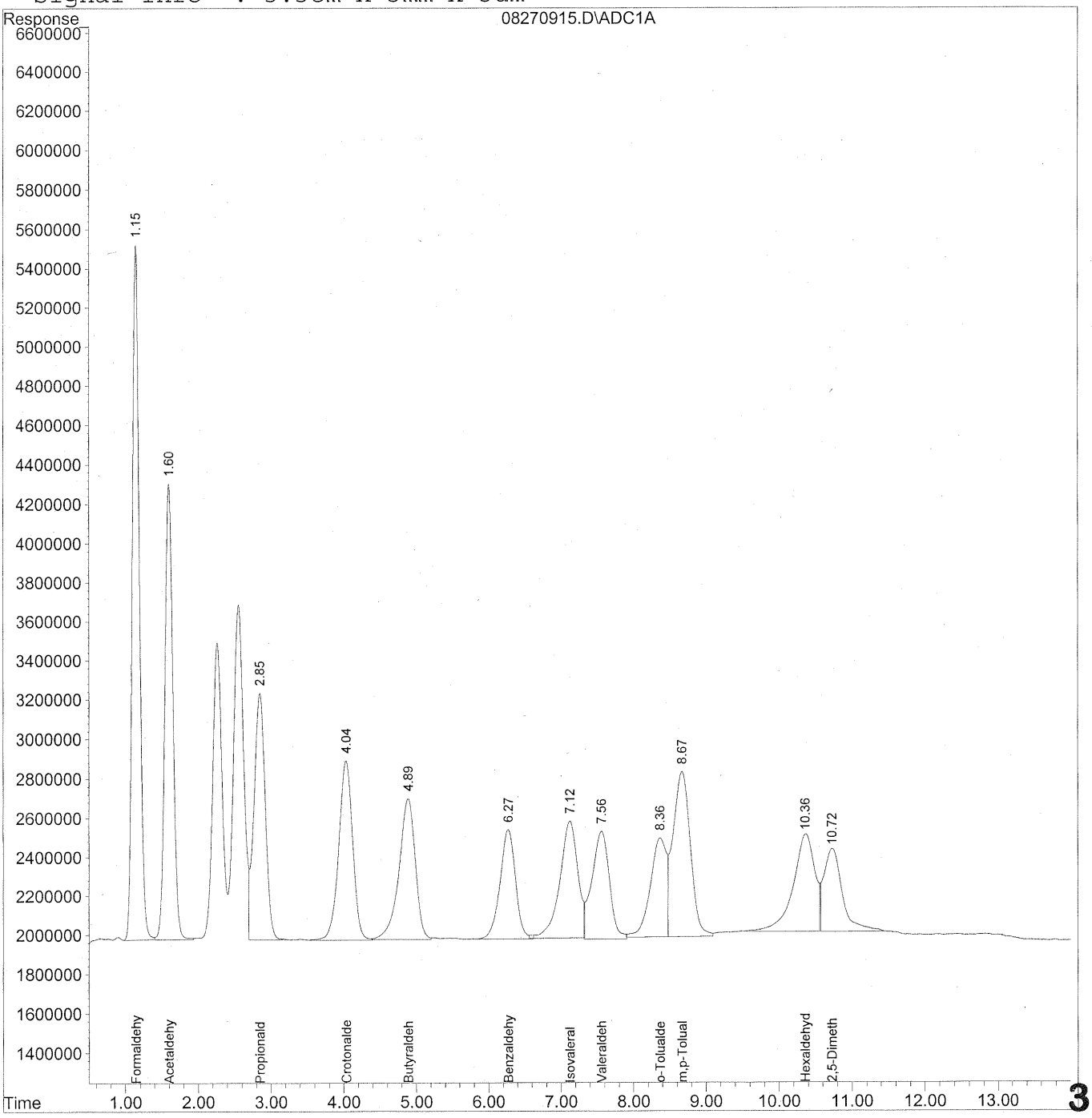
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.14	258608681	1408.687	ng/ml
2)	Acetaldehyde	1.59	194077889	1384.061	ng/ml
3)	Propionaldehyde	2.82	146239778	1370.631	ng/ml
4)	Crotonaldehyde	3.98	133040312	1365.704	ng/ml
5)	Butyraldehyde	4.83	123280292	1395.582	ng/ml
6)	Benzaldehyde	6.22	91701289	1392.170	ng/ml
7)	Isovaleraldehyde	7.09	110300508	1409.573	ng/ml
8)	Valeraldehyde	7.54	95980847	1305.773	ng/ml
9)	o-Tolualdehyde	8.35	82388557	1412.688	ng/ml
10)	m,p-Tolualdehyde	8.67	148951031	2758.587	ng/ml
11)	Hexaldehyde	10.43	94972826	1410.269	ng/ml
12)	2,5-Dimethylbenzaldehyde	10.80	63578554	1297.167	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:52 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 : Sat Aug 29 17:49:00 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\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:52 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 : Sat Aug 29 17:49:00 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

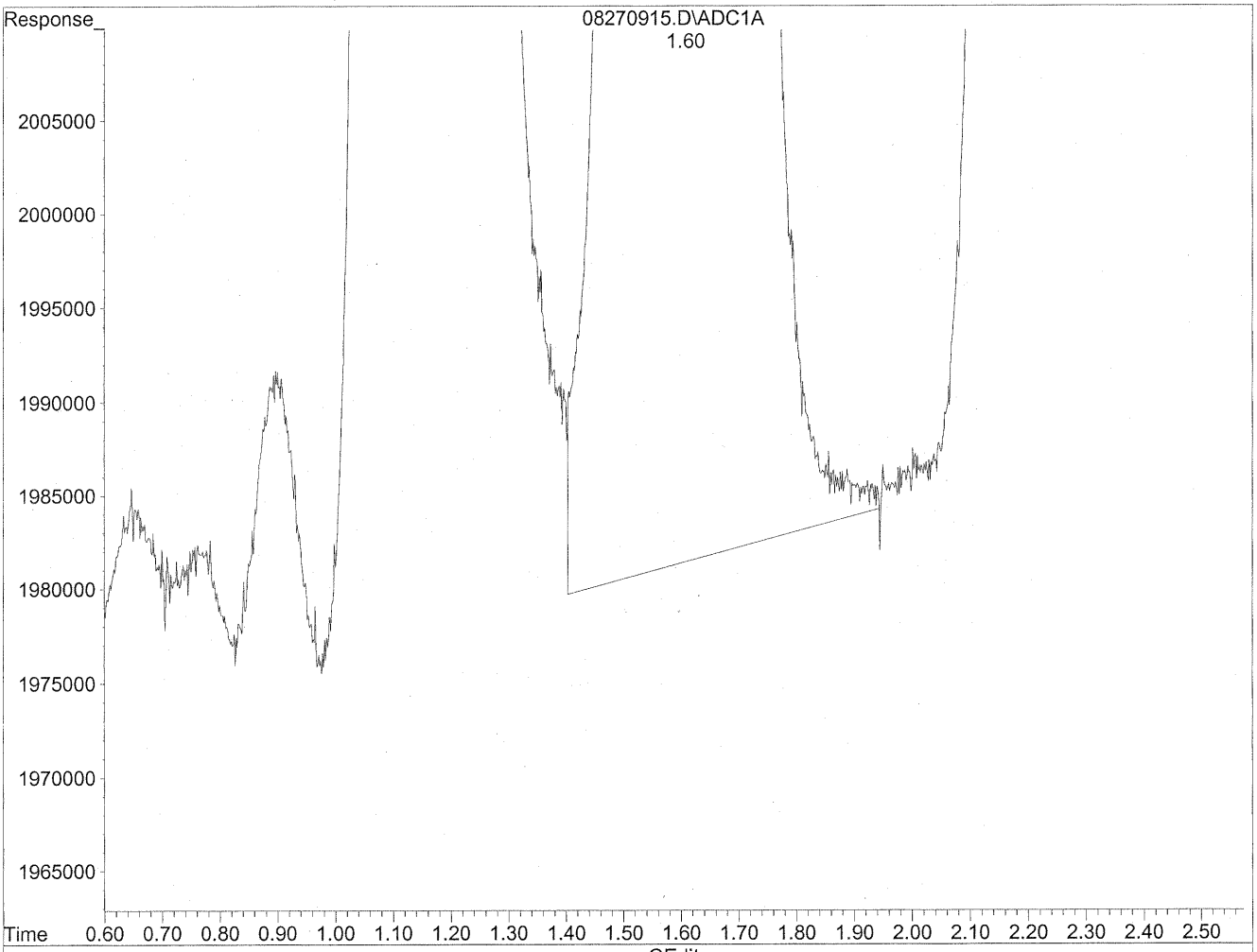
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.15	236705173	1289.375	ng/ml
2)	Acetaldehyde	1.60	179688294	1281.442	ng/mlm
3)	Propionaldehyde	2.85	136845423	1282.582	ng/mlm
4)	Crotonaldehyde	4.04	122298348	1255.434	ng/mlm
5)	Butyraldehyde	4.89	113329905	1282.940	ng/mlm
6)	Benzaldehyde	6.27	84558301	1283.728	ng/mlm
7)	Isovaleraldehyde	7.12	104917810	1340.786	ng/ml
8)	Valeraldehyde	7.56	93919272	1277.726	ng/mlm
9)	o-Tolualdehyde	8.37	78916199	1353.149	ng/ml
10)	m,p-Tolualdehyde	8.67	142355511	2636.438	ng/ml
11)	Hexaldehyde	10.36	102492195	1521.926	ng/mlm
12)	2,5-Dimethylbenzaldehyde	10.72	74449913	1518.971	ng/mlm

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 13:06 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.60min 1272.083ng/ml

response 178375946

315

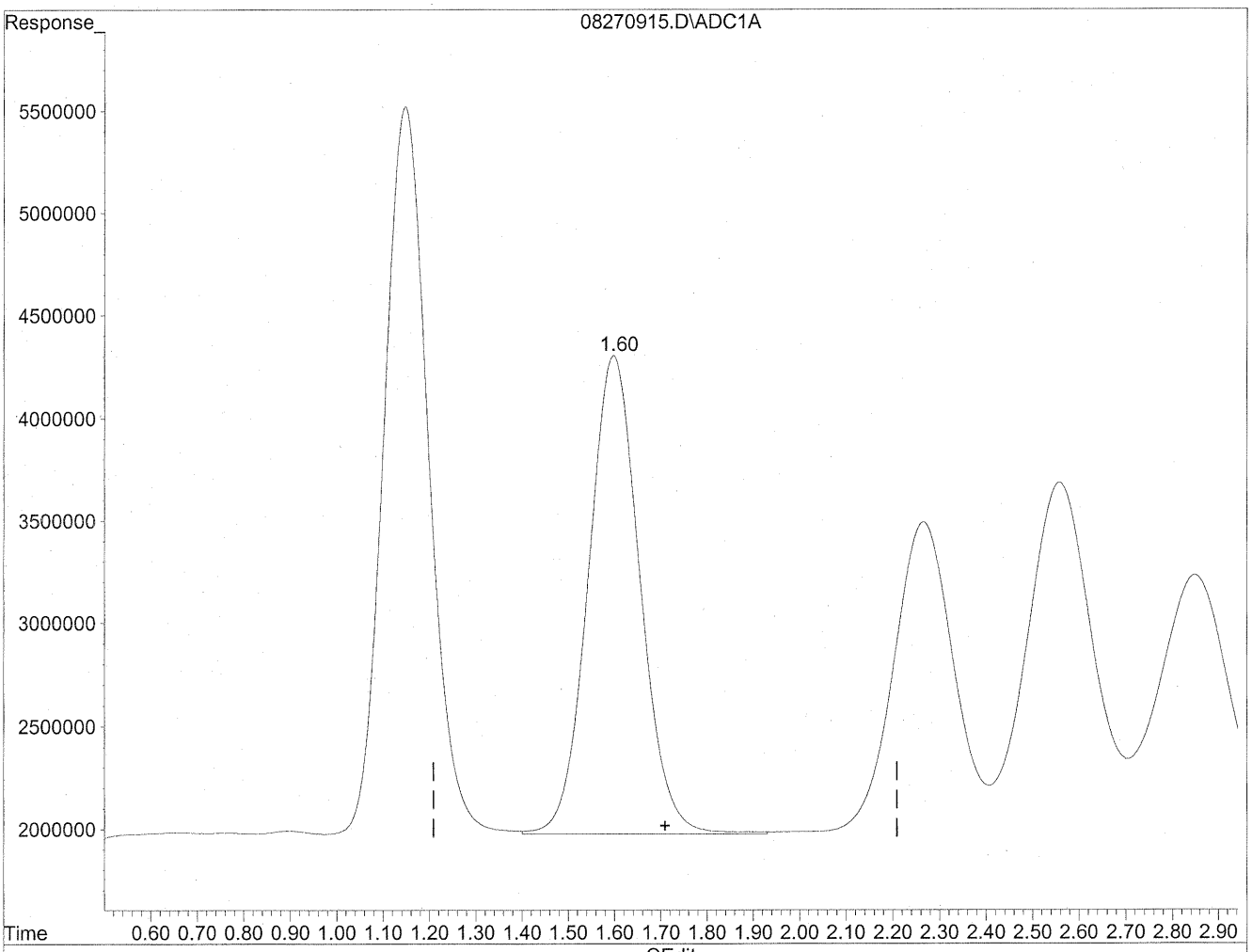
(+) = Expected Retention Time

08270915.D TO110709.M Sat Aug 29 17:44:03 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 27 13:06 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 : Sat Aug 29 16:33:38 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.60min 1285.302ng/ml m

response 180229578

HC 8/31/09 BC

8/31/09

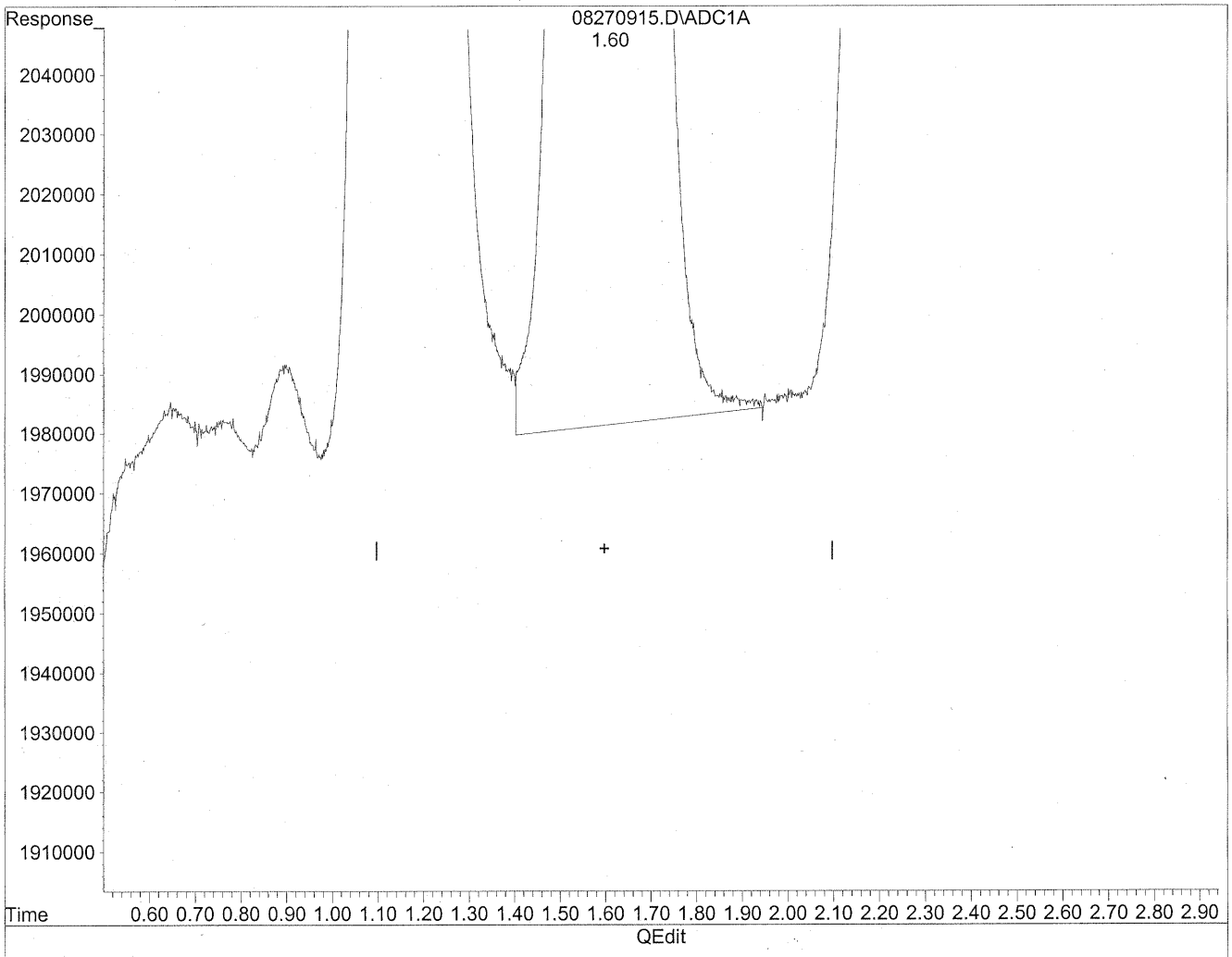
316

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.60min 1272.083ng/ml

response 178375946

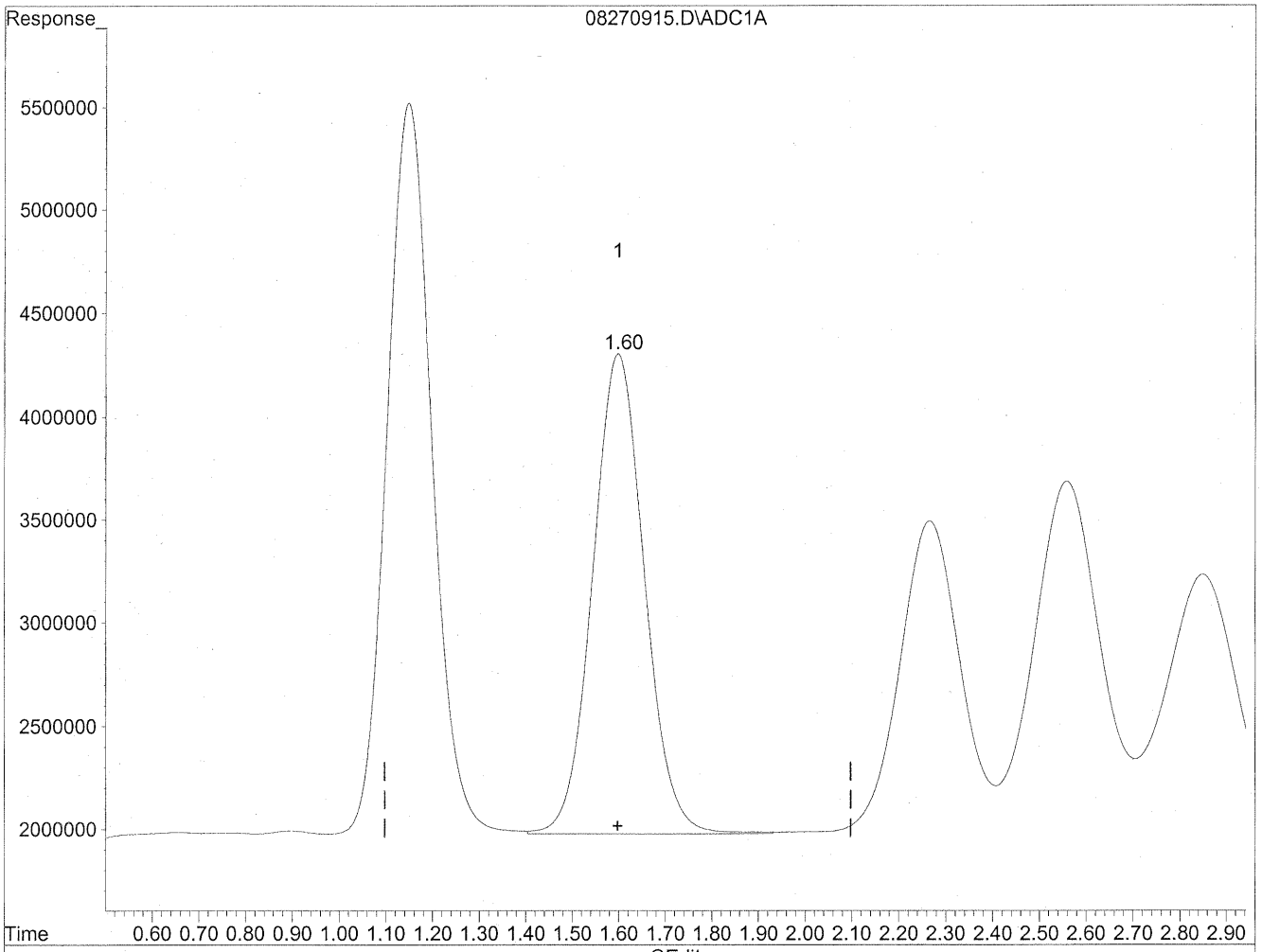
317

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.60min 1281.442ng/ml m

response 179688294

HC 8/30/09
BC

8/31/09

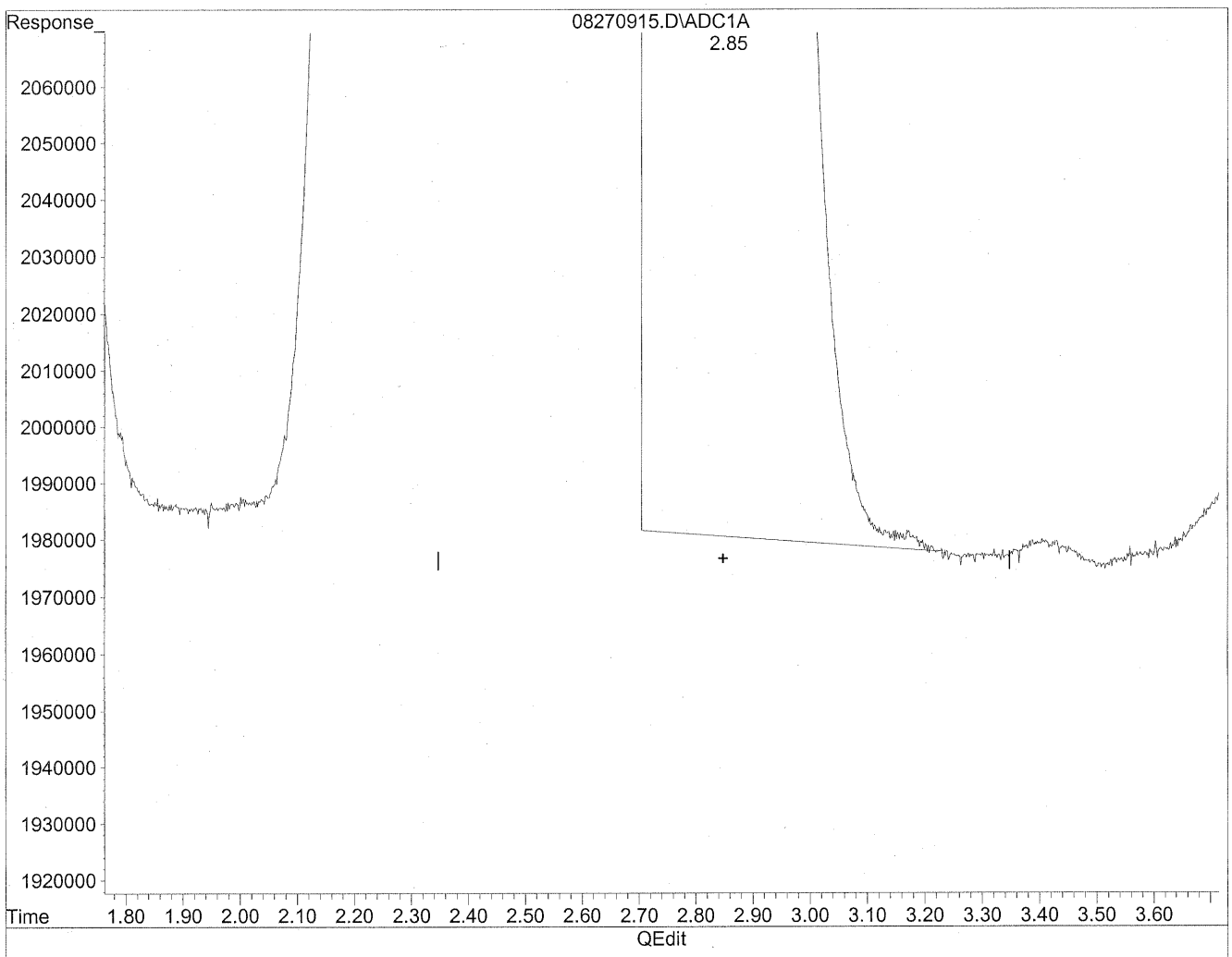
318

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.85min 1257.093ng/ml

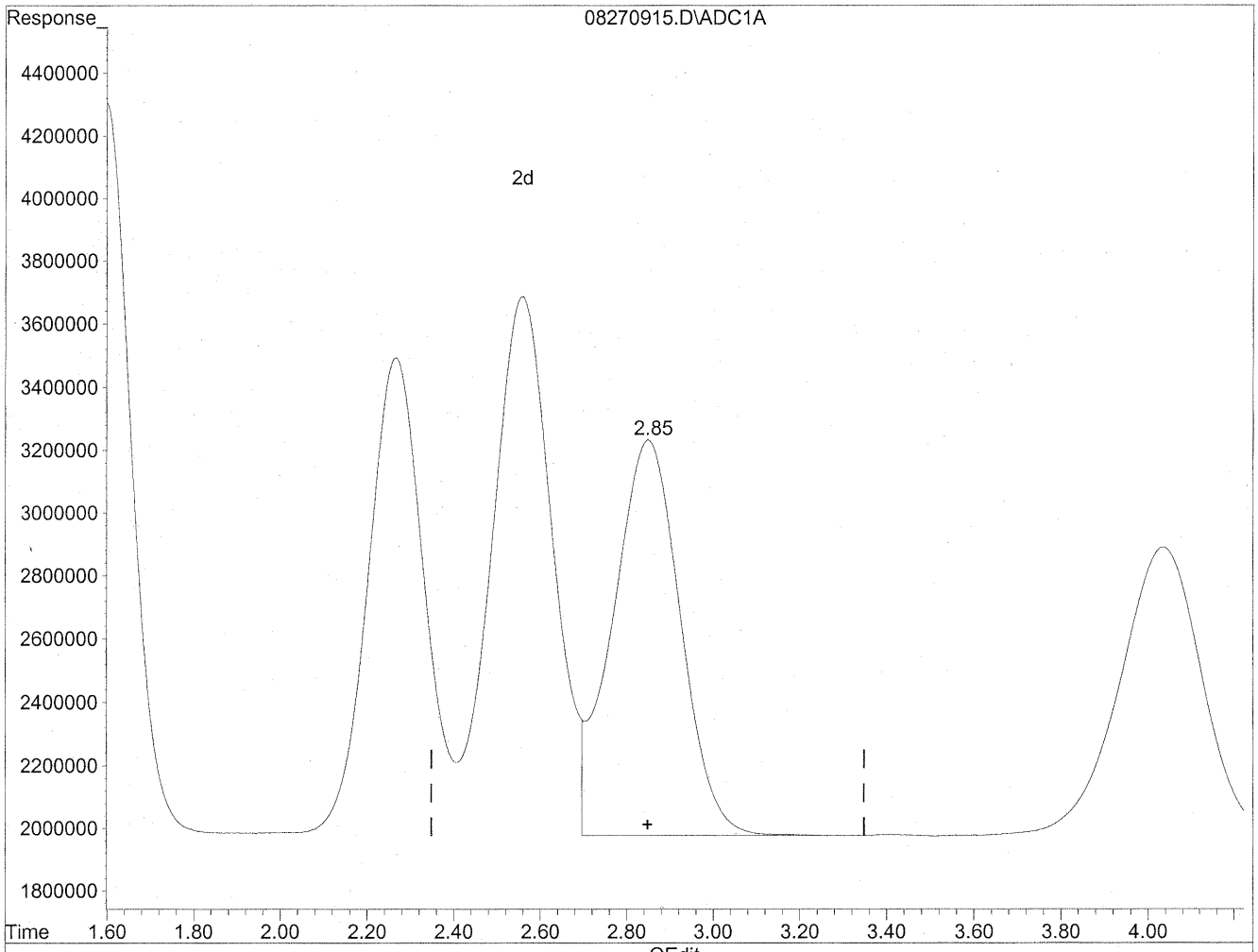
response 134125861

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.85min 1282.582ng/ml m

response 136845423

JC 8/31/09 BC

Wog 8/31/09

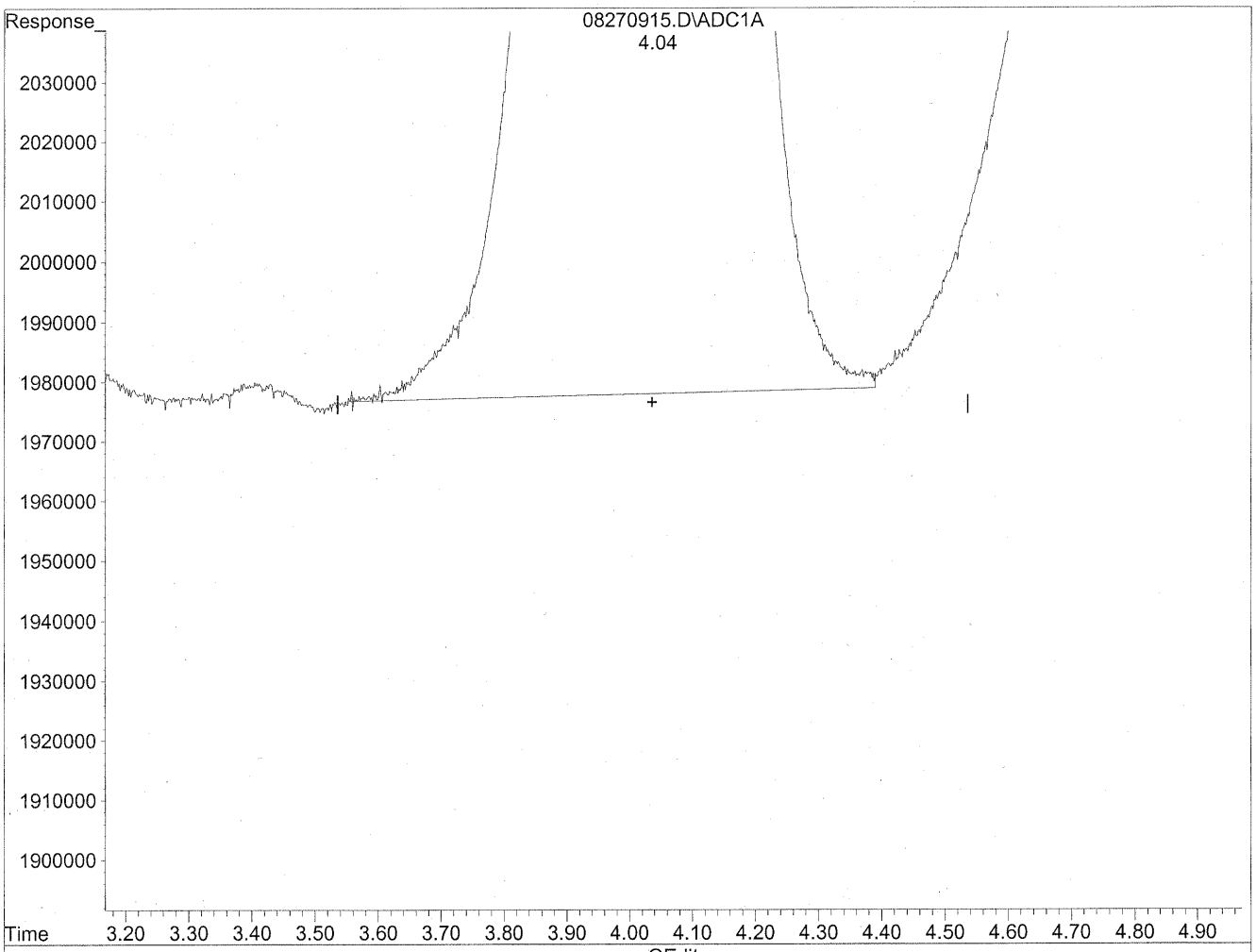
320

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.03min 1238.377ng/ml

response 120636689

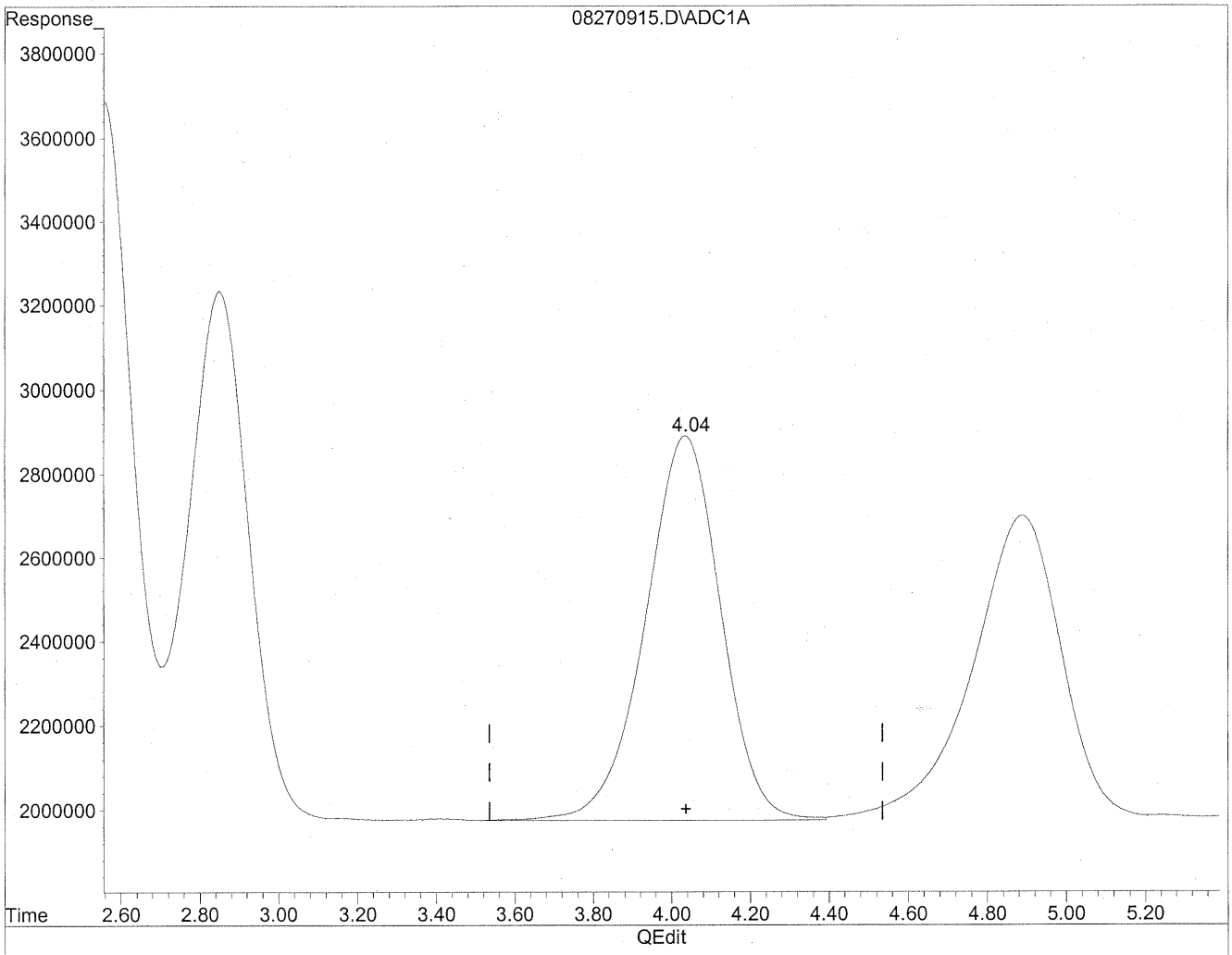
321

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.04min 1255.434ng/ml m

response 122298348

HC 8/30/09
BC

WH 8/31/09

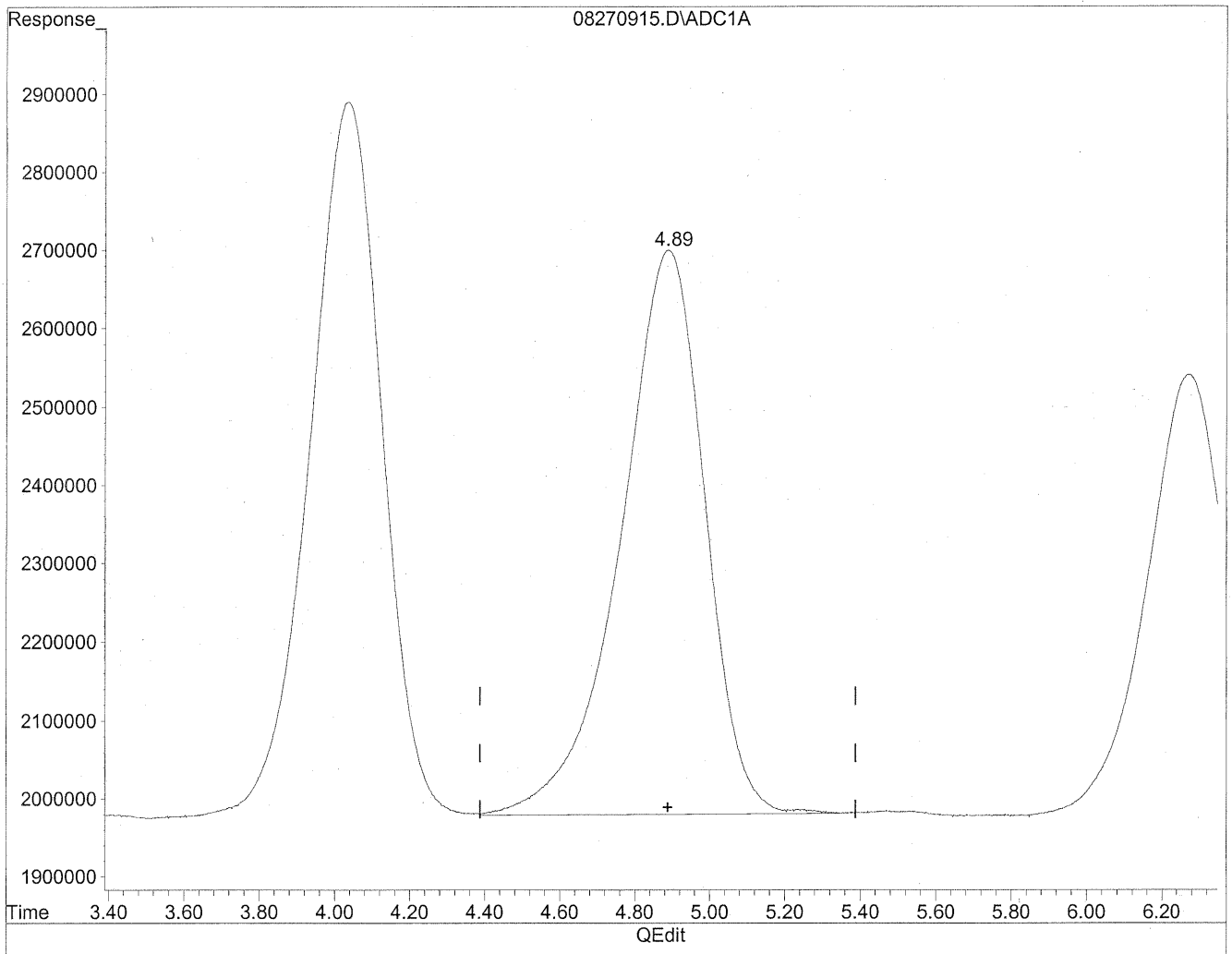
322

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.89min 1265.793ng/ml

response 111815270

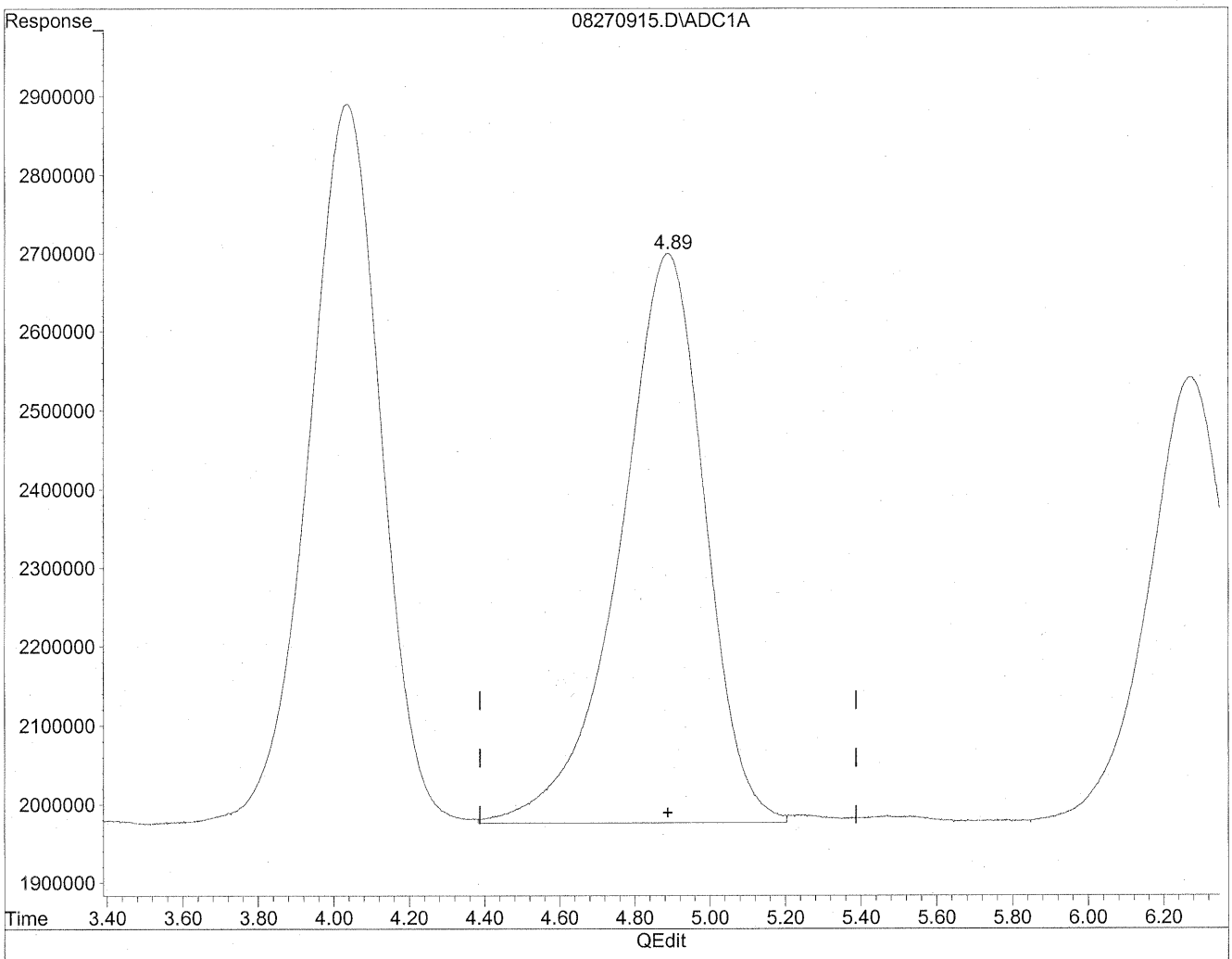
(+) = Expected Retention Time

08270915.D TO110709.M Sat Aug 29 17:50:28 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.89min 1282.940ng/ml m

response 113329905

HC 8/30/09 BC

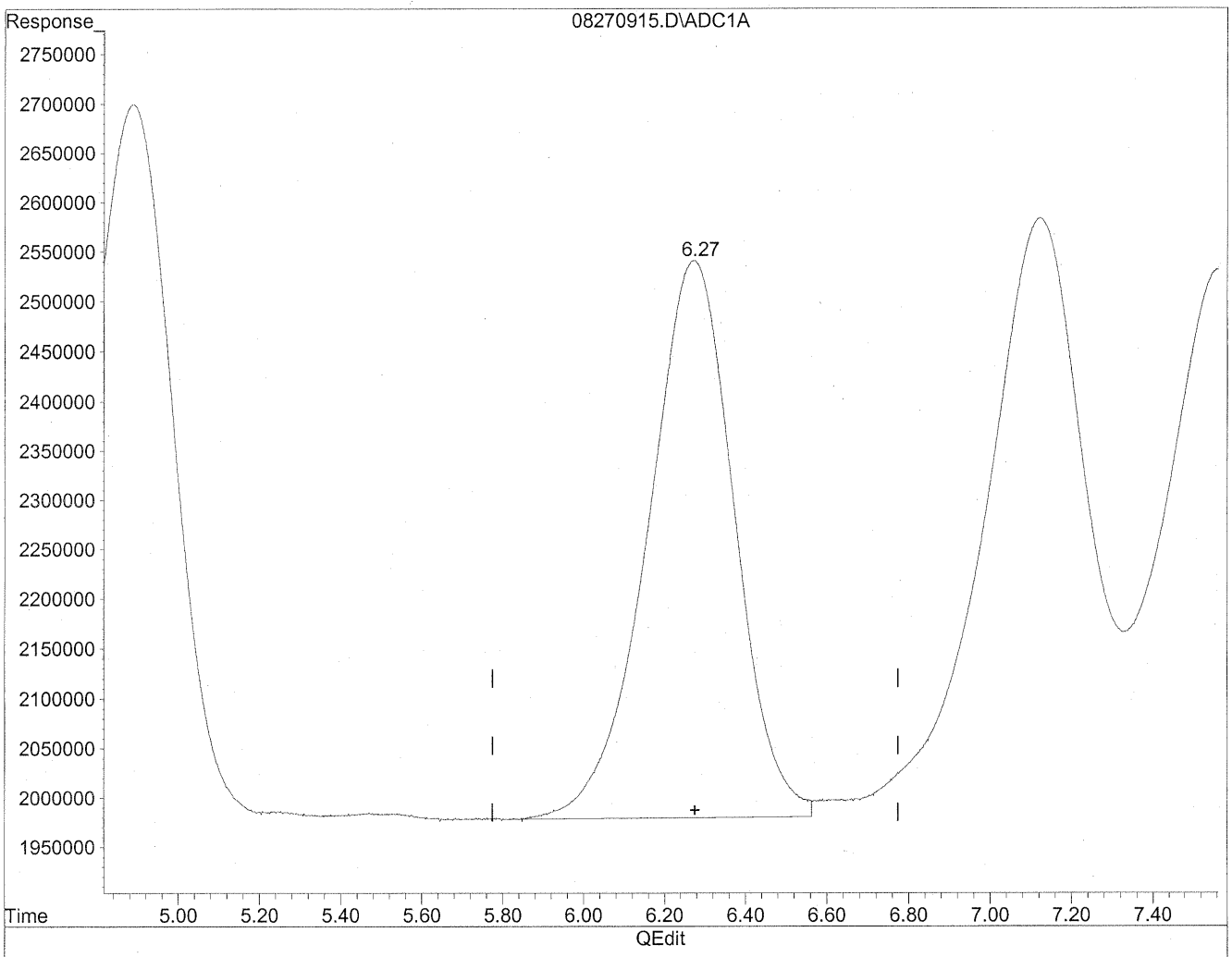
LA 8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.27min 1263.121ng/ml

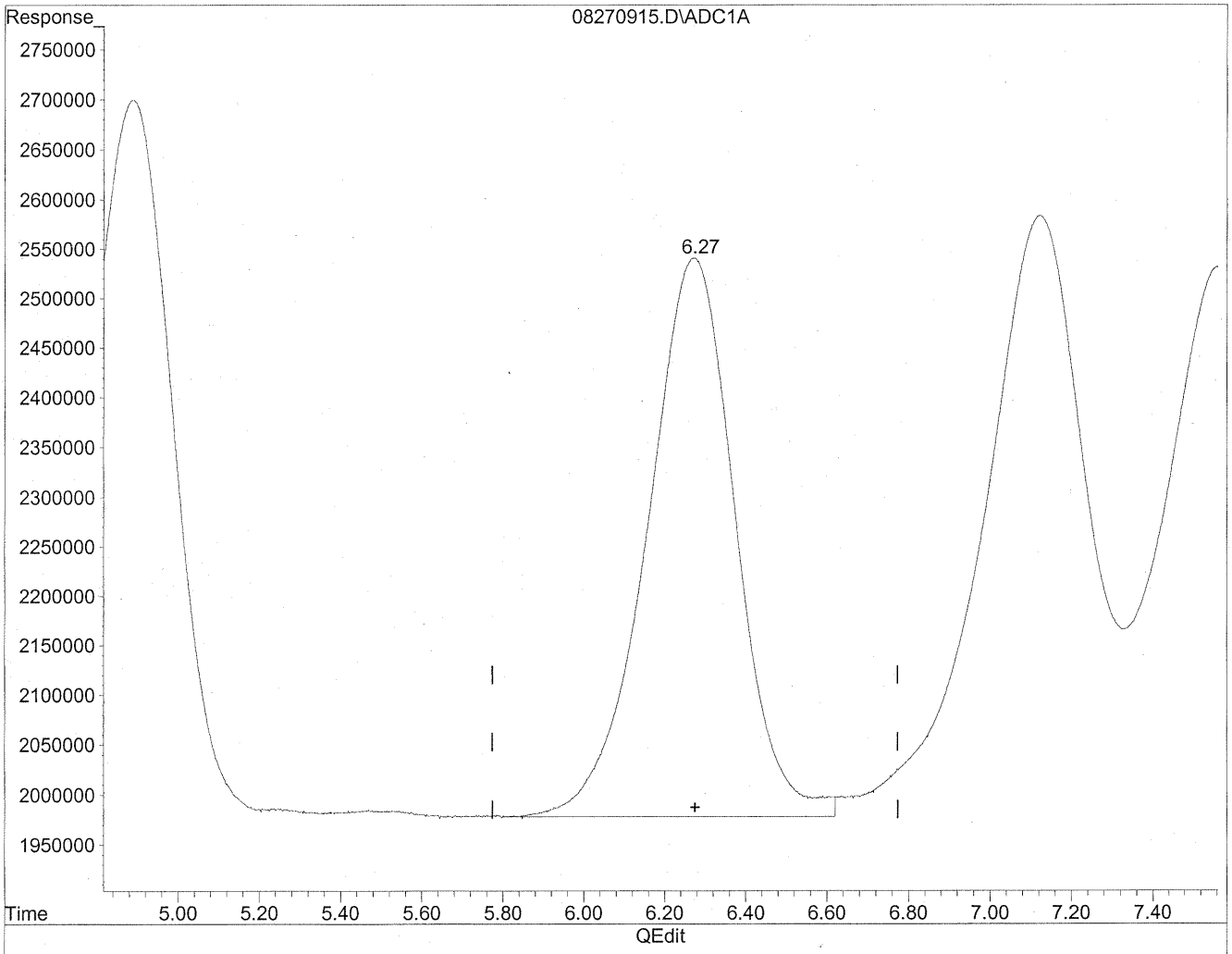
response 83200901

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.27min 1283.728ng/ml m

response 84558301

HC 8/30/09

BC

8/31/09

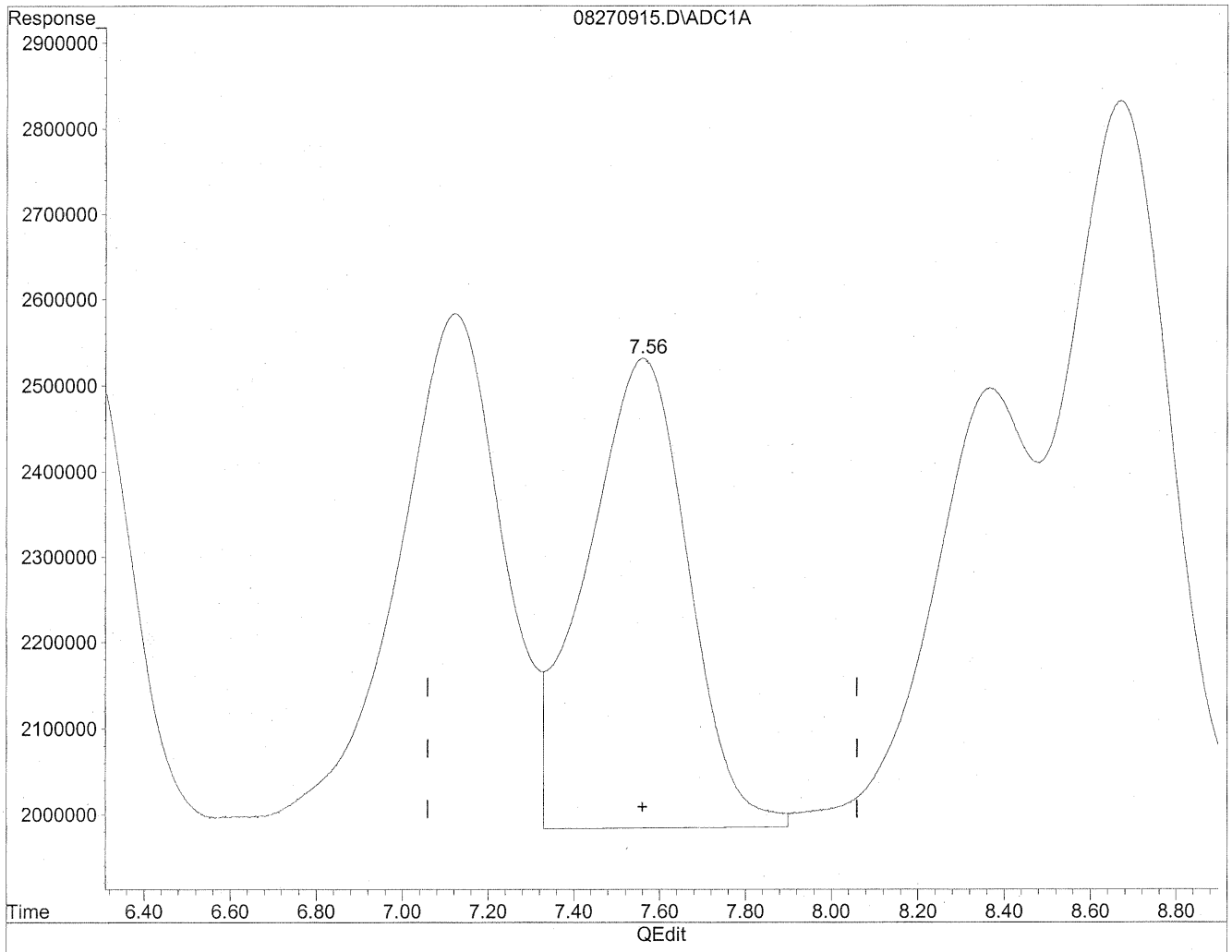
326

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.56min 1225.744ng/ml

response 90098322

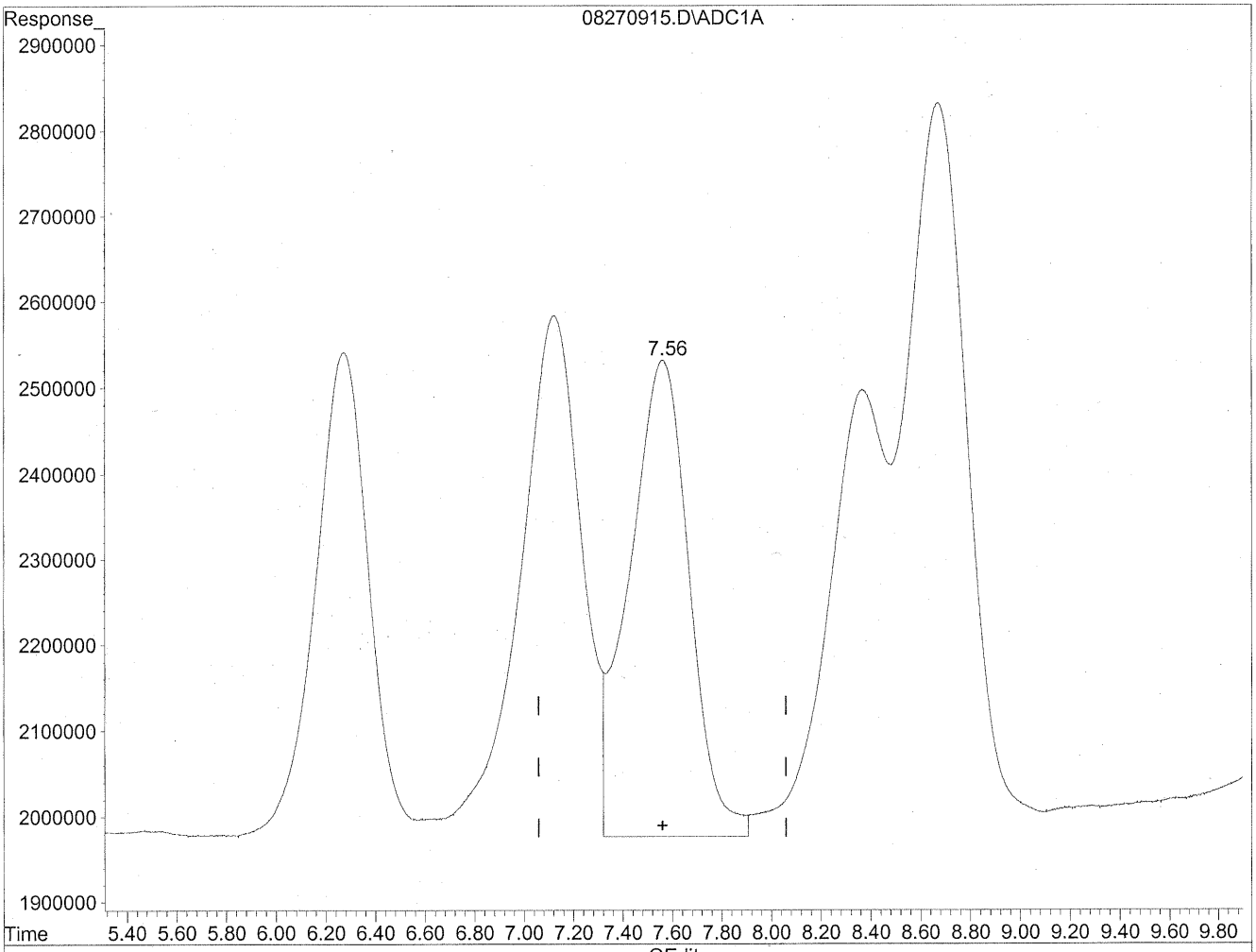
(+) = Expected Retention Time

08270915.D TO110709.M Sat Aug 29 17:51:03 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.56min 1277.726ng/ml m

response 93919272

HC 8/30/07
BL

WB 8/31/07

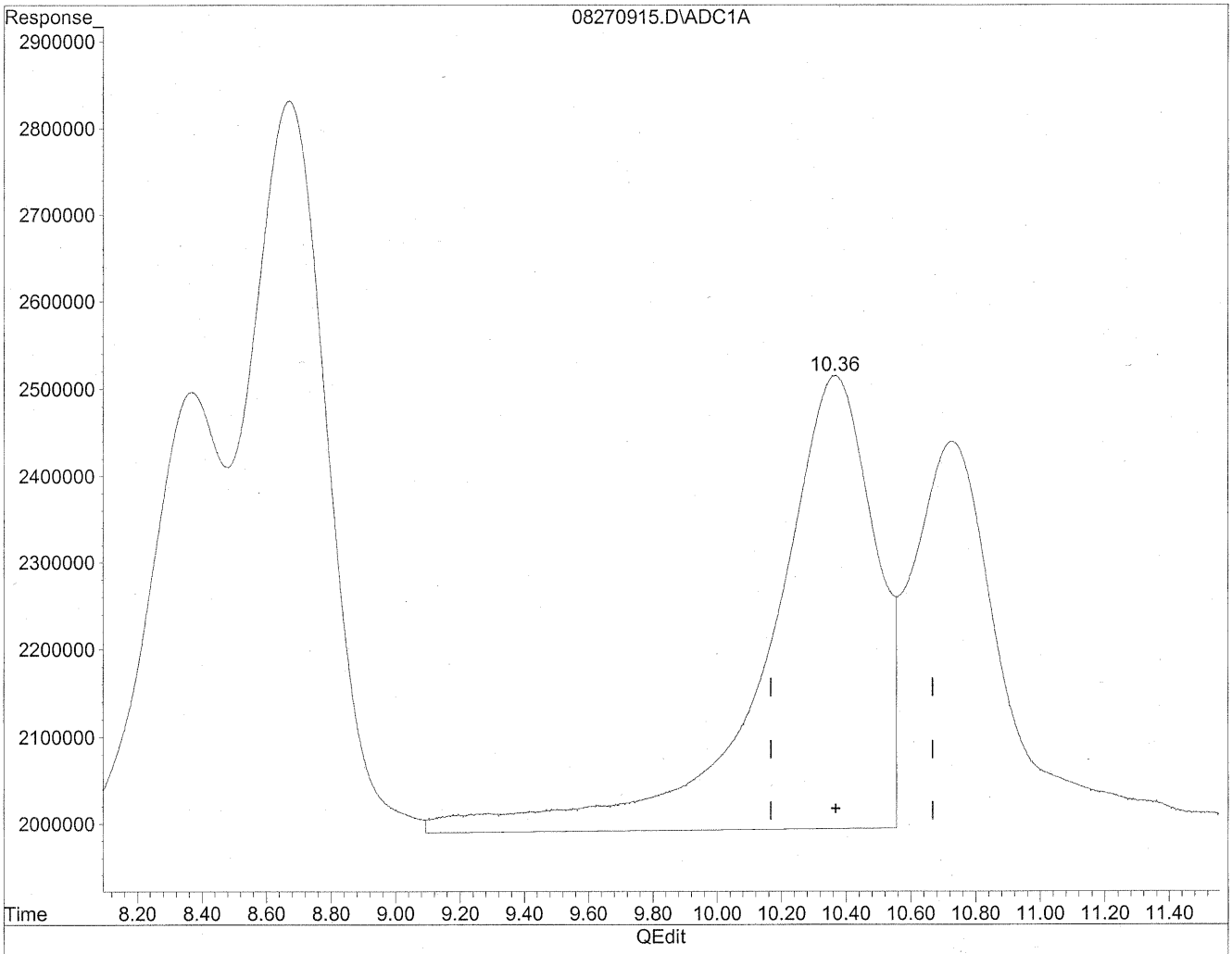
328

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.37min 1794.915ng/ml

response 120876298

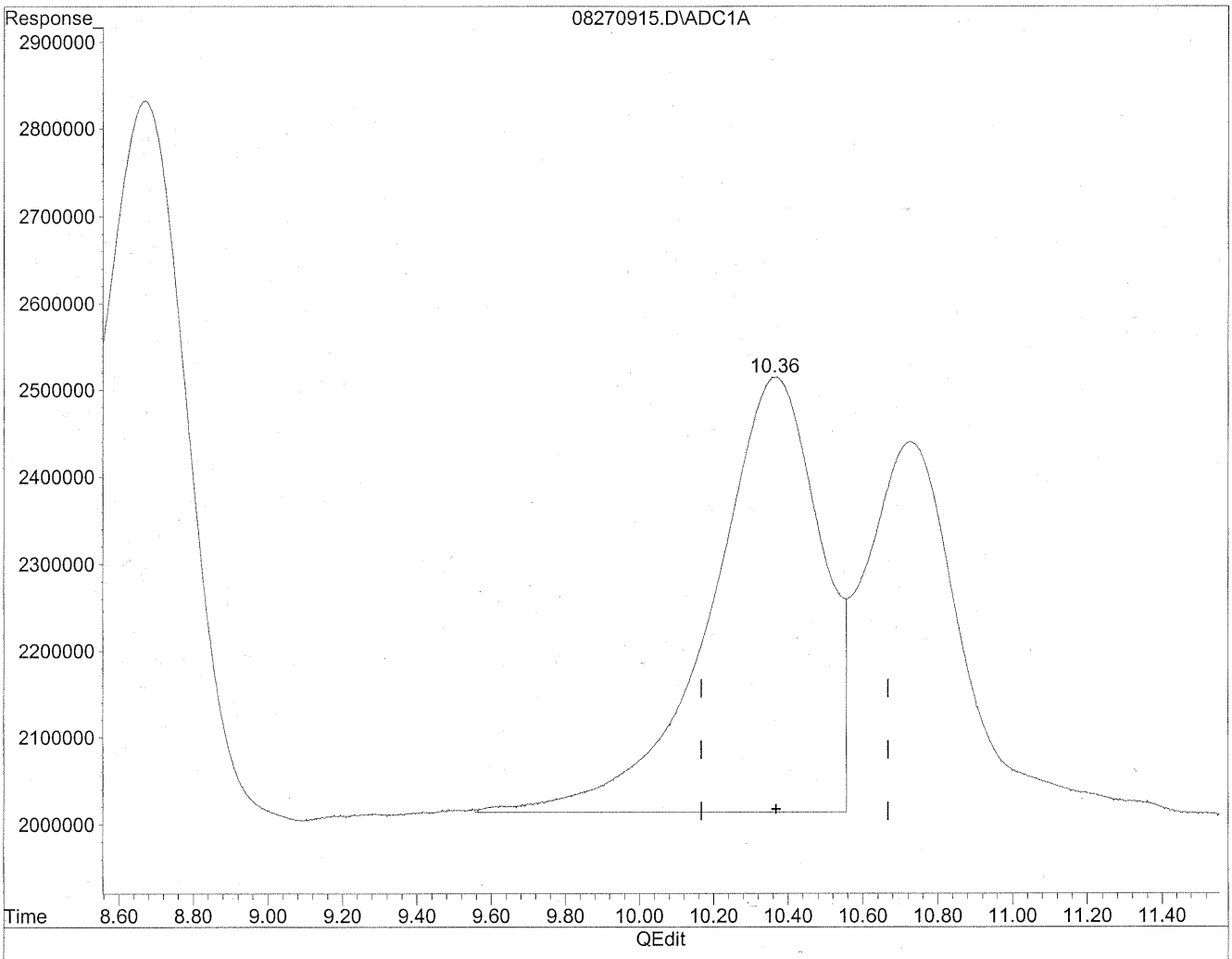
329

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\27\08270915.D Vial: 15
Acq On : 27 Aug 2009 12:36 pm Operator: HC
Sample : CCV 1500ng/ml S21-08270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 17:49 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 17:49:00 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.36min 1521.926ng/ml m

response 102492195

HC 8/30/09
BC

330

(+) = Expected Retention Time

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

Injection Log

Directory: j:\lc01\data\to11\2009_08\26

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	1	08260901.d	1.	prime		26 Aug 109 12::
2	2	08260902.d	1.	1500ng/ml TO11A std S21-08250901		26 Aug 109 12::
3	3	08260903.d	1.	ACN blank Lot CY023		26 Aug 109 12::
4	4	08260904.d	1.	MB front lot 5855/5994 1.0ml		26 Aug 109 12::
5	5	08260905.d	1.	MB back lot 5855/5994 1.0ml		26 Aug 109 12::
6	6	08260906.d	1.	P0902942-001 back 1.0ml		26 Aug 109 12::
7	7	08260907.d	1.	P0902942-002 back 1.0ml		26 Aug 109 12::
8	8	08260908.d	1.	P0902942-003 back 1.0ml		26 Aug 109 12::
9	9	08260909.d	1.	P0902942-004 back 1.0ml		26 Aug 109 12::
10	10	08260910.d	1.	P0902942-005 back 1.0ml		26 Aug 109 12::
11	11	08260911.d	1.	P0902942-006 back 1.0ml		26 Aug 109 12::
12	12	08260912.d	1.	P0902946-001 back 1.0ml		26 Aug 109 12::
13	13	08260913.d	1.	P0902946-002 back 1.0ml		26 Aug 109 12::
14	14	08260914.d	1.	P0902946-003 back 1.0ml		26 Aug 109 12::
15	15	08260915.d	1.	P0902946-004 back 1.0ml		26 Aug 109 12::
16	16	08260916.d	1.	CCV 1500ng/ml S21-08250901		26 Aug 109 12::
17	17	08260917.d	1.	P0902946-005 back 1.0ml		26 Aug 109 12::
18	17	08260918.d	1.	P0902946-005dup back 1.0ml		26 Aug 109 12::
19	18	08260919.d	1.	P0902946-006 back 1.0ml		26 Aug 109 12::
20	19	08260920.d	1.	P0902946-007 back 1.0ml		26 Aug 109 12::
21	20	08260921.d	1.	P0902946-008 back 1.0ml		26 Aug 109 13::
22	21	08260922.d	1.	P0902946-009 back 1.0ml		26 Aug 109 13::
23	22	08260923.d	1.	P0902946-0010 back 1.0ml		26 Aug 109 13::
24	23	08260924.d	1.	P0902946-0011 back 1.0ml		26 Aug 109 13::
25	24	08260925.d	1.	P0902946-012 back 1.0ml		26 Aug 109 13::
26	25	08260926.d	1.	P0902946-013 back 1.0ml		26 Aug 109 13::
27	26	08260927.d	1.	CCV 1500ng/ml S21-08250901		26 Aug 109 13::
28	27	08260928.d	1.	ACN lot blk CY023		26 Aug 109 13::
29	28	08260929.d	1.	MB front lot 5855/5994 1.0ml		27 Aug 109 13::
30	29	08260930.d	1.	MB back lot 5855/5994 1.0ml		27 Aug 109 13::
31	30	08260931.d	1.	P0902946-015 back 1.0ml		27 Aug 109 13::
32	31	08260932.d	1.	P0902946-016 back 1.0ml		27 Aug 109 13::
33	32	08260933.d	1.	P0902946-017 back 1.0ml		27 Aug 109 12::
34	33	08260934.d	1.	P0902946-018 back 1.0ml		27 Aug 109 12::
35	34	08260935.d	1.	P0902946-019 back 1.0ml		27 Aug 109 12::
36	35	08260936.d	1.	P0902946-020 back 1.0ml		27 Aug 109 12::
37	36	08260937.d	1.	P0902946-021 back 1.0ml		27 Aug 109 12::
38	37	08260938.d	1.	P0902946-022 back 1.0ml		27 Aug 109 12::
39	38	08260939.d	1.	P0902946-023 back 1.0ml		27 Aug 109 12::
40	39	08260940.d	1.	CCV 1500ng/ml S21-08250901		27 Aug 109 12::
41	40	08260941.d	1.	P0902946-024 back 1.0ml		27 Aug 109 12::
42	41	08260942.d	1.	P0902946-025 back 1.0ml		27 Aug 109 12::
43	42	08260943.d	1.	P0902946-026 back 1.0ml		27 Aug 109 12::
44	43	08260944.d	1.	P0902946-027 back 1.0ml		27 Aug 109 12::
45	44	08260945.d	1.	P0902946-028 back 1.0ml		27 Aug 109 12::
46	45	08260946.d	1.	P0902946-029 back 1.0ml		27 Aug 109 12::
47	46	08260947.d	1.	P0902946-014 back 1.0ml		27 Aug 109 12::
48	47	08260948.d	1.	P0902942-001 front 1.0ml		27 Aug 109 12::
49	47	08260949.d	1.	P0902942-001dup front 1.0ml		27 Aug 109 12::
50	48	08260950.d	1.	ACN wash		27 Aug 109 12::
51	49	08260951.d	1.	CCV 1500ng/ml S21-08250901		27 Aug 109 12::

333

Injection Log

Directory: j:\lc01\data\to11\2009_08\27

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	1	08270901.d	1.	1500ng/ml TO11A std S21-08270903		27 Aug 109 12::
2	2	08270902.d	1.	ACN blank Lot CY023		27 Aug 109 12::
3	3	08270903.d	1.	MB front lot 5855/5994 1.0ml		27 Aug 109 12::
4	4	08270904.d	1.	MB back lot 5855/5994 1.0ml		27 Aug 109 12::
5	5	08270905.d	1.	P0902942-002 front 1.0ml		27 Aug 109 13::
6	6	08270906.d	1.	P0902942-003 front 1.0ml		27 Aug 109 13::
7	7	08270907.d	1.	P0902942-004 front 1.0ml		27 Aug 109 13::
8	8	08270908.d	1.	P0902942-005 front 1.0ml		27 Aug 109 13::
9	9	08270909.d	1.	P0902942-006 front 1.0ml		27 Aug 109 13::
10	10	08270910.d	1.	P0902946-001 front 1.0ml		27 Aug 109 13::
11	11	08270911.d	1.	P0902946-002 front 1.0ml		27 Aug 109 13::
12	12	08270912.d	1.	P0902946-003 front 1.0ml		27 Aug 109 13::
13	13	08270913.d	1.	P0902946-004 front 1.0ml		27 Aug 109 13::
14	14	08270914.d	1.	ACN wash		27 Aug 109 13::
15	15	08270915.d	1.	CCV 1500ng/ml S21-08270903		27 Aug 109 13::
16	16	08270916.d	1.	P0902946-005 front 1.0ml		27 Aug 109 13::
17	17	08270917.d	1.	P0902946-006 front 1.0ml		27 Aug 109 12::
18	17	08270918.d	1.	P0902946-006dup front 1.0ml		27 Aug 109 12::
19	18	08270919.d	1.	P0902946-007 front 1.0ml		27 Aug 109 12::
20	19	08270920.d	1.	P0902946-008 front 1.0ml		27 Aug 109 12::
21	20	08270921.d	1.	P0902946-009 front 1.0ml		27 Aug 109 12::
22	21	08270922.d	1.	P0902946-010 front 1.0ml		27 Aug 109 12::
23	22	08270923.d	1.	P0902946-011 front 1.0ml		27 Aug 109 12::
24	23	08270924.d	1.	P0902946-012 front 1.0ml		27 Aug 109 12::
25	24	08270925.d	1.	P0902946-013 front 1.0ml		27 Aug 109 12::
26	25	08270926.d	1.	P0902946-014 front 1.0ml		27 Aug 109 12::
27	26	08270927.d	1.	ACN wash		27 Aug 109 12::
28	27	08270928.d	1.	CCV 1500ng/ml S21-08270903		27 Aug 109 12::
29	28	08270929.d	1.	ACN blk lot CY023		27 Aug 109 12::
30	29	08270930.d	1.	MB front lot 5855/5994 1.0ml		27 Aug 109 12::
31	30	08270931.d	1.	MB back lot 5855/5994 1.0ml		27 Aug 109 12::
32	31	08270932.d	1.	P0902946-015 front 1.0ml		27 Aug 109 12::
33	32	08270933.d	1.	P0902946-016 front 1.0ml		27 Aug 109 12::
34	33	08270934.d	1.	P0902946-017 front 1.0ml		27 Aug 109 12::
35	34	08270935.d	1.	P0902946-018 front 1.0ml		27 Aug 109 12::
36	35	08270936.d	1.	P0902946-019 front 1.0ml		27 Aug 109 12::
37	36	08270937.d	1.	P0902946-020 front 1.0ml		27 Aug 109 12::
38	37	08270938.d	1.	P0902946-021 front 1.0ml		27 Aug 109 12::
39	38	08270939.d	1.	P0902946-022 front 1.0ml		27 Aug 109 12::
40	39	08270940.d	1.	P0902946-023 front 1.0ml		27 Aug 109 12::
41	40	08270941.d	1.	P0902946-024 front 1.0ml		27 Aug 109 12::
42	41	08270942.d	1.	ACN wash		27 Aug 109 12::
43	42	08270943.d	1.	CCV 1500ng/ml S21-08270903		27 Aug 109 12::
44	43	08270944.d	1.	P0902946-025 front 1.0ml		27 Aug 109 12::
45	43	08270945.d	1.	P0902946-025dup front 1.0ml		27 Aug 109 12::
46	44	08270946.d	1.	P0902946-026 front 1.0ml		27 Aug 109 12::
47	45	08270947.d	1.	P0902946-027 front 1.0ml		27 Aug 109 12::
48	46	08270948.d	1.	P0902946-028 front 1.0ml		27 Aug 109 12::
49	47	08270949.d	1.	P0902946-029 front 1.0ml		27 Aug 109 12::
50	48	08270950.d	1.	ACN wash		27 Aug 109 12::
51	49	08270951.d	1.	CCV 1500ng/ml S21-08270903		27 Aug 109 12::
52	50	08270952.d	1.	ACN blk lot CY023		27 Aug 109 12::
53	51	08270953.d	1.	MB front lot 5855/5994 1.0ml		27 Aug 109 13::
54	52	08270954.d	1.	MB back lot 5855/5994 1.0ml		27 Aug 109 13::
55	53	08270955.d	1.	P0902964-001 back1.0ml		27 Aug 109 13::
56	54	08270956.d	1.	P0902964-002 back 1.0ml		27 Aug 109 13::
57	55	08270957.d	1.	P0902964-003 back 1.0ml		27 Aug 109 13::

Injection Log

Directory: j:\lc01\data\to11\2009_08\27

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
58	56	08270958.d	1.	P0902964-004 back 1.0ml		27 Aug 109 13::
59	57	08270959.d	1.	P0902964-005 back 1.0ml		27 Aug 109 13::
60	58	08270960.d	1.	P0902964-006 back 1.0ml		27 Aug 109 13::
61	59	08270961.d	1.	P0902965-001 back 1.0ml		28 Aug 109 13::
62	60	08270962.d	1.	P0902965-002 back 1.0ml		28 Aug 109 13::
63	61	08270963.d	1.	P0902965-003 back 1.0ml		28 Aug 109 13::
64	62	08270964.d	1.	P0902965-004 back 1.0ml		28 Aug 109 13::
65	63	08270965.d	1.	ACN wash		28 Aug 109 12::
66	64	08270966.d	1.	CCV 1500ng/ml S21-08270903		28 Aug 109 12::
67	65	08270967.d	1.	P0902965-005 back 1.0ml		28 Aug 109 12::
68	65	08270968.d	1.	P0902965-005dup back 1.0ml		28 Aug 109 12::
69	66	08270969.d	1.	P0902965-006 back 1.0ml		28 Aug 109 12::
70	67	08270970.d	1.	P0902965-007 back 1.0ml		28 Aug 109 12::
71	68	08270971.d	1.	P0902965-008 back 1.0ml		28 Aug 109 12::
72	69	08270972.d	1.	P0902965-009 back 1.0ml		28 Aug 109 12::
73	70	08270973.d	1.	P0902965-010 back 1.0ml		28 Aug 109 12::
74	71	08270974.d	1.	P0902965-011 back 1.0ml		28 Aug 109 12::
75	72	08270975.d	1.	P0902965-012 back 1.0ml		28 Aug 109 12::
76	73	08270976.d	1.	ACN wash		28 Aug 109 12::
77	74	08270977.d	1.	CCV 1500ng/ml S21-08270903		28 Aug 109 12::