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

LABORATORY REPORT

September 16, 2009

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

RE: 16512

Dear Brian:

Enclosed are the results of the samples submitted to our laboratory on August 21, 2009. For your reference, these analyses have been assigned our service request number P0902912.

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 257 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, Incorporated CAS Project No: P0902912
Project: 16512

CASE NARRATIVE

The samples were received intact under chain of custody on August 21, 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 results of analyses are given in the attached laboratory report. All results are intended to be considered in their entirety, and Columbia Analytical Services, Inc. (CAS) is not responsible for utilization of less than the complete report.

Client: Environmental Health & Engineering, Incorporate
Project: 16512

Service Request: P0902912

SAMPLE CROSS-REFERENCE

SAMPLE #	CLIENT SAMPLE ID	DATE	TIME
P0902912-001	102509	8/20/09	00:00
P0902912-002	102511	8/20/09	00:00
P0902912-003	102512	8/20/09	00:00
P0902912-004	102513	8/20/09	00:00
P0902912-005	102514	8/20/09	00:00

3

Environmental Health & Engineering, Inc.

CHAIN OF CUSTODY FORM

DATE: 20 AUG 2009

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

90902912

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)	102509	AIR	EPA TO-11	102.0 L
(2)	102511	"	FULL LIST	101.5 "
(3)	102512	"	"	108.68 "
(4)	102513	"	"	102.96 "
(5)	102514	"	"	

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 mfrazala@eheinc.com

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

Relinquished by: _____ of Environmental Health & Engineering, Inc. Date: 8/20/09

Received by: _____ of (company name) CAS Date: 8/21/09 0940

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, Incorporated Work order: P0902912

Project: 16512

Sample(s) received on: 08/21/09 Date opened: 08/21/09 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 15 °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
P0902912-001.01	Silica Gel DNPH Tube					
P0902912-002.01	Silica Gel DNPH Tube					
P0902912-003.01	Silica Gel DNPH Tube					
P0902912-004.01	Silica Gel DNPH Tube					
P0902912-005.01	Silica Gel DNPH Tube					

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

Sample -001 was received with the endcaps detached.

*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: **102509**
Client Project ID: **16512**

CAS Project ID: P0902912
CAS Sample ID: P0902912-001

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: 8/20/09
Date Received: 8/21/09
Date Analyzed: 8/25/09
Desorption Volume: 1.0 ml
Volume Sampled: 102 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**5,600**	**55**	0.98	**45**	0.80	
75-07-0	**Acetaldehyde**	**6,000**	**59**	0.98	**33**	0.54	**BT**
123-38-6	**Propionaldehyde**	**570**	**5.6**	0.98	**2.4**	0.41	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.98	ND	0.34	
123-72-8	**Butyraldehyde**	**590**	**5.8**	0.98	**2.0**	0.33	
100-52-7	**Benzaldehyde**	**400**	**3.9**	0.98	**0.90**	0.23	
590-86-3	**Isovaleraldehyde**	**230**	**2.2**	0.98	**0.64**	0.28	
110-62-3	**Valeraldehyde**	**1,500**	**14**	0.98	**4.1**	0.28	
529-20-4	o-Tolualdehyde	< 100	ND	0.98	ND	0.20	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	2.0	ND	0.40	
66-25-1	**n-Hexaldehyde**	**6,500**	**63**	0.98	**15**	0.24	
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	ND	0.98	ND	0.18	

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

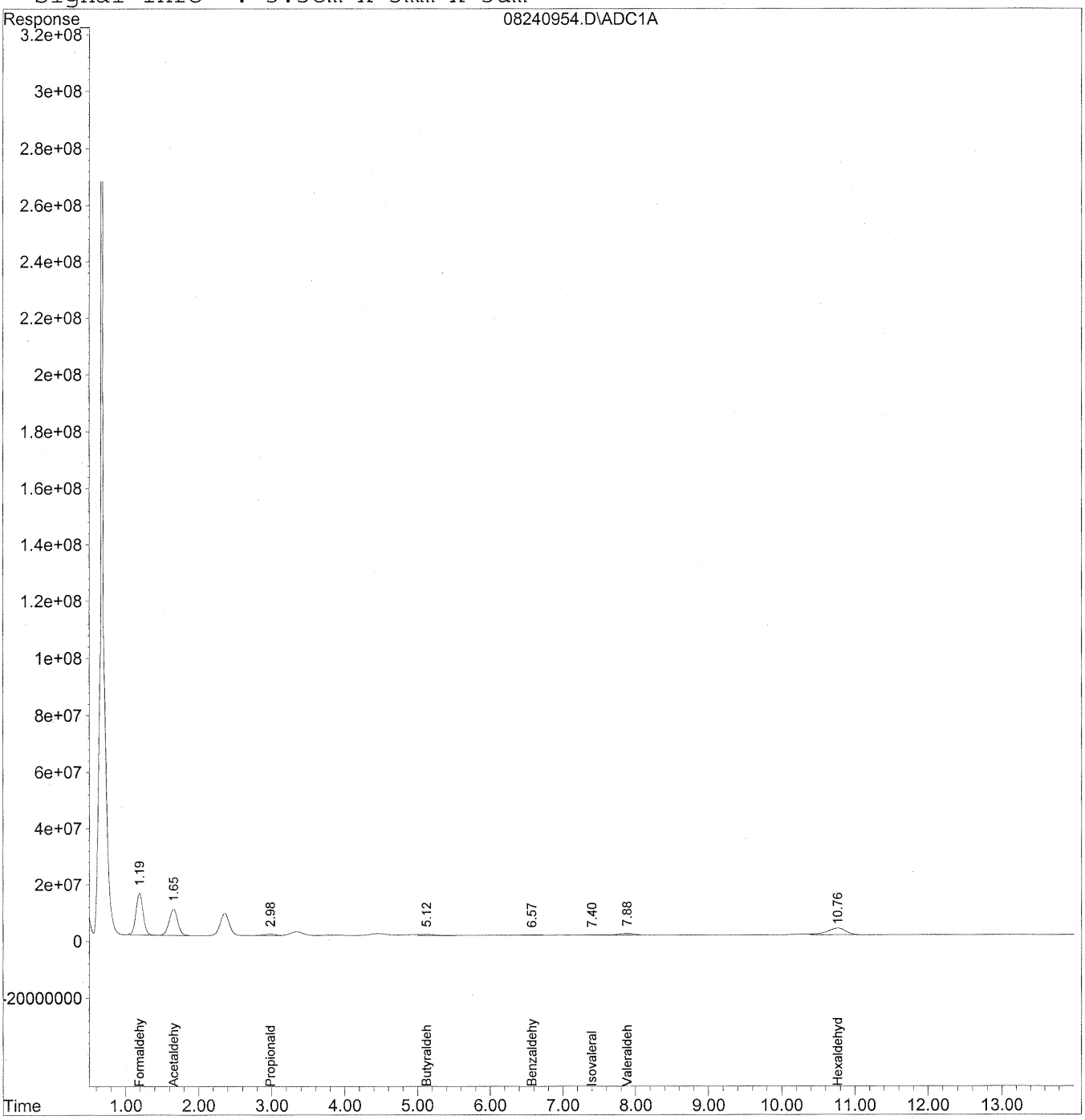
Verified By: RG Date: 9/2/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 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 Aug 25 10:21:57 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)

Data File : J:\LC01\DATA\TO11\2009_08\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 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 Aug 25 10:21:57 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

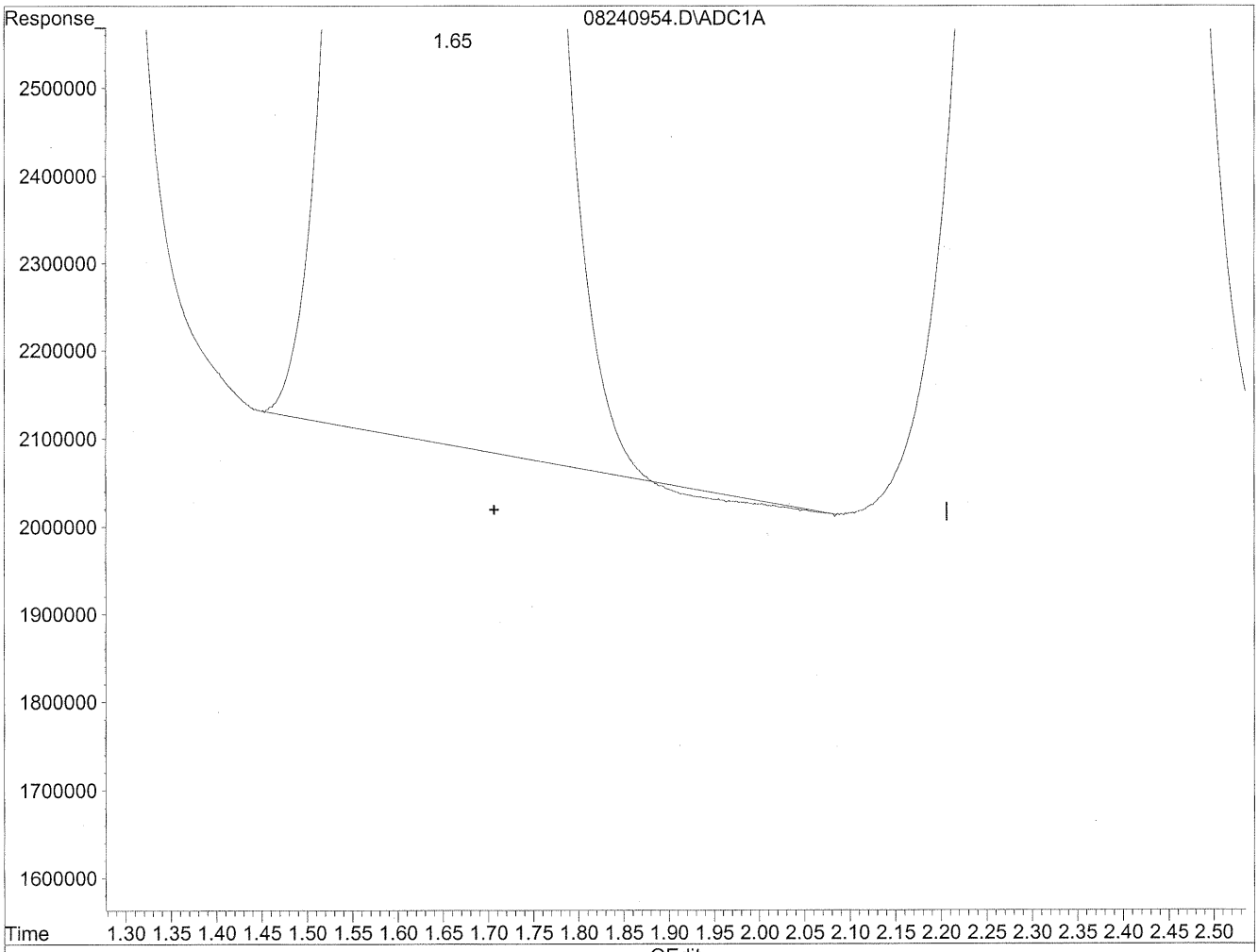
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.19	996681423	5429.098	ng/ml
2)	Acetaldehyde	1.65	766485914	5466.173	ng/mlm
3)	Propionaldehyde	2.98	60828006	570.110	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	5.12	51896644	587.491	ng/mlm
6)	Benzaldehyde	6.57	26127092	396.650	ng/mlm
7)	Isovaleraldehyde	7.40	17869572	228.362	ng/mlm
8)	Valeraldehyde	7.88	107912963	1468.104	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	10.76f	434454506	6451.295	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\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:56 19109 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

1.65min 5456.761ng/ml

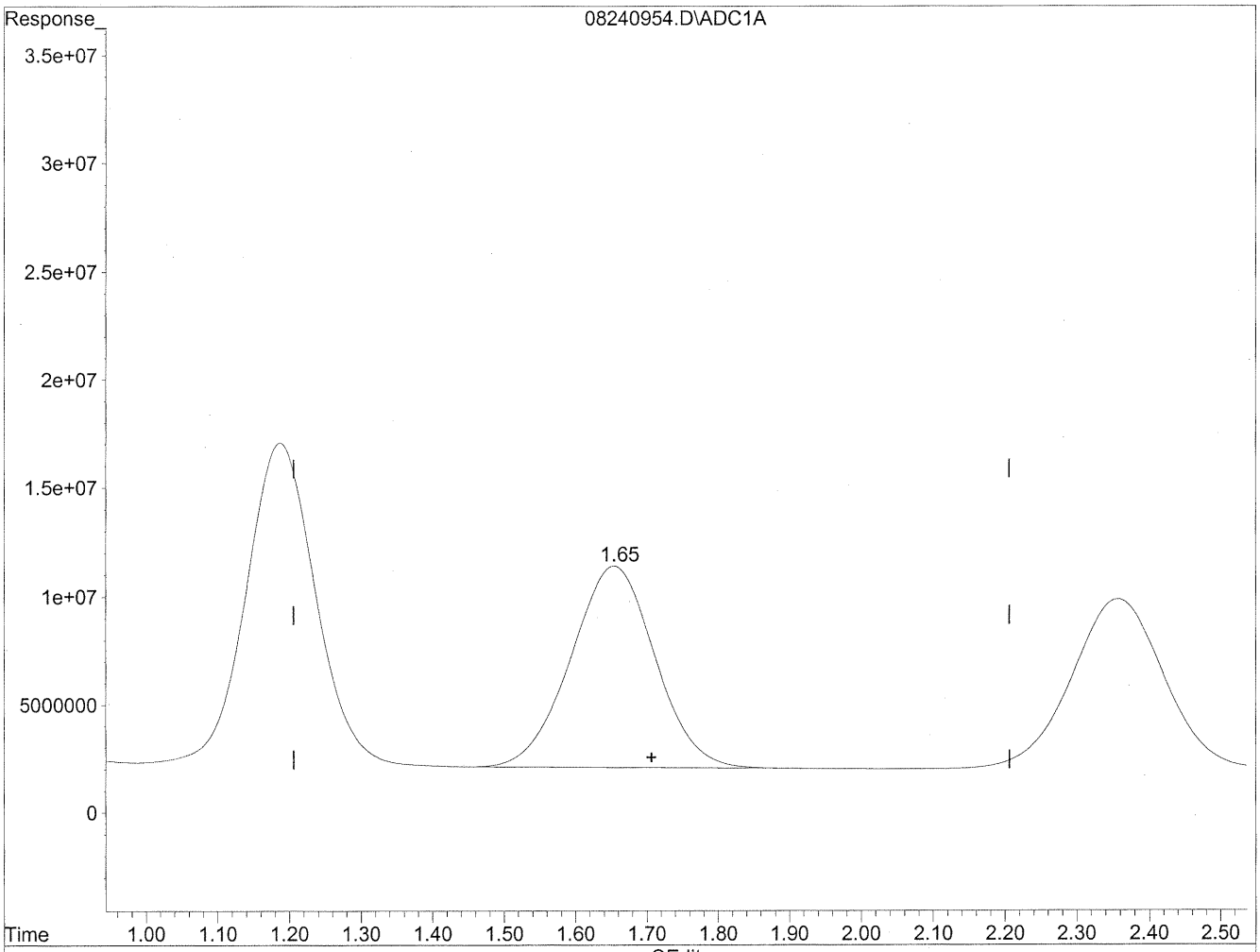
response 765166188

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:56 19109 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

1.65min 5466.173ng/ml m

response 766485914

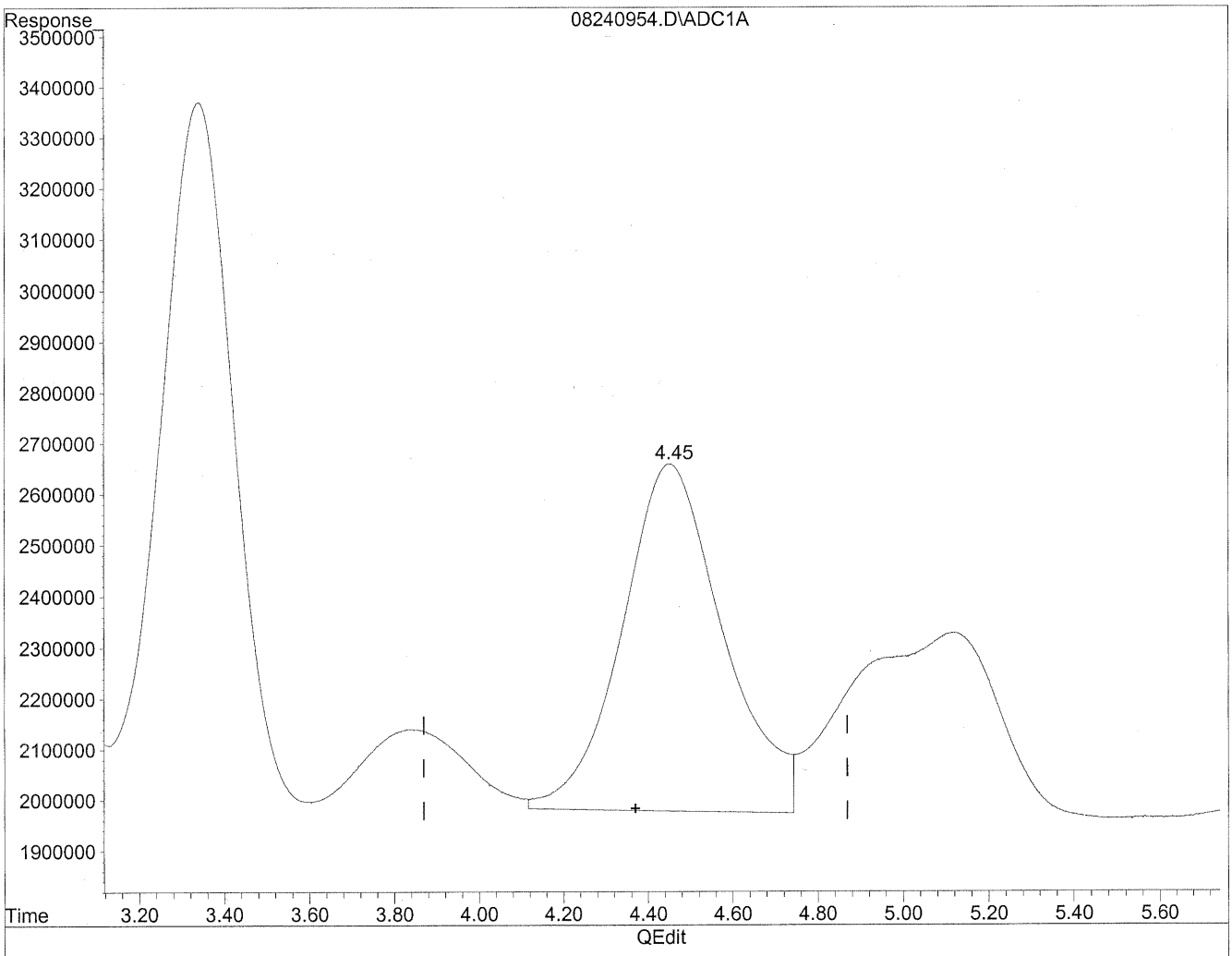
HC 8/27/09 LC

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:56 19109 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



(4) Crotonaldehyde

4.45min 1160.408ng/ml

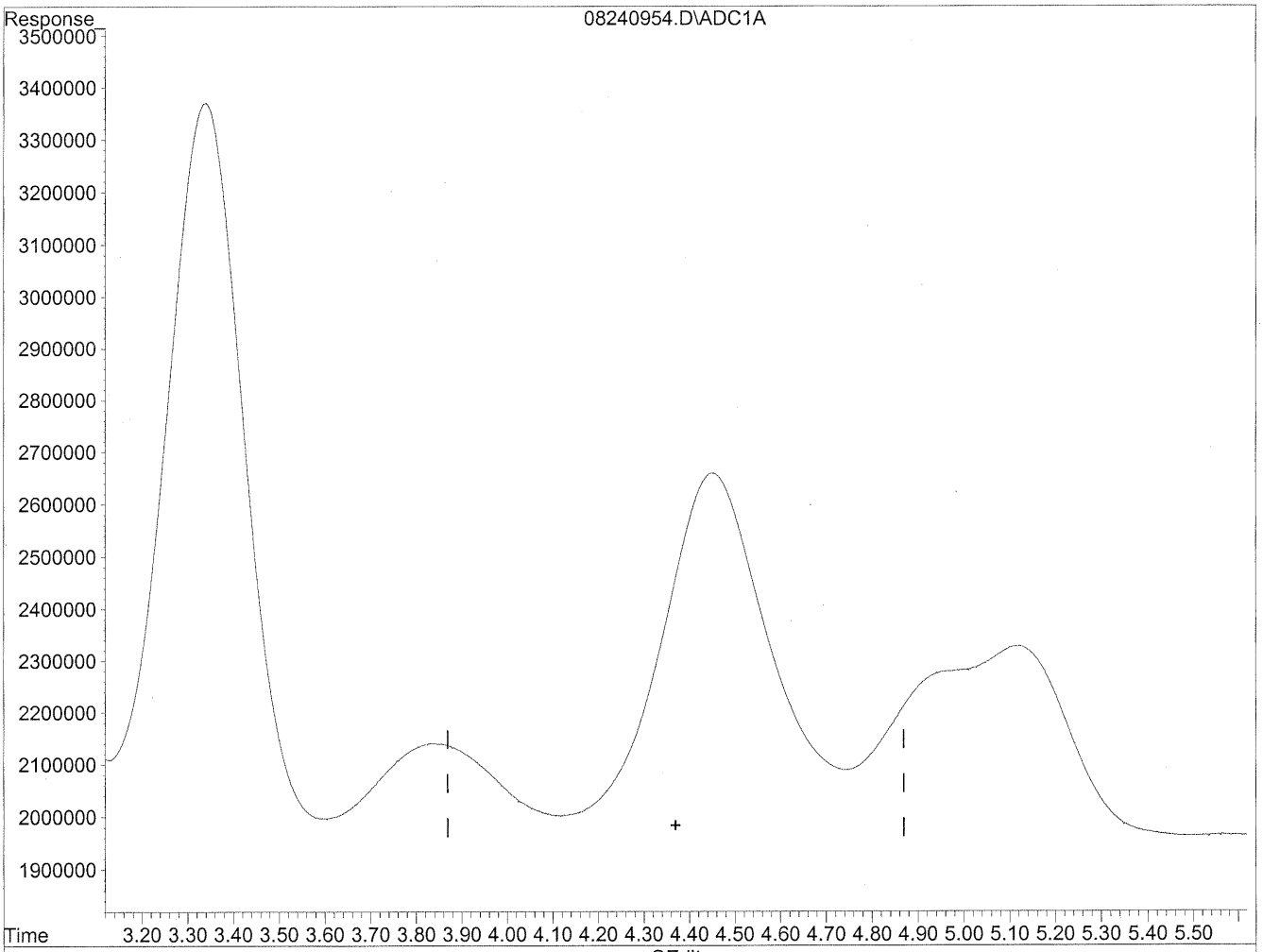
response 113041394

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:56 19109 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



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/24/09 WP

13

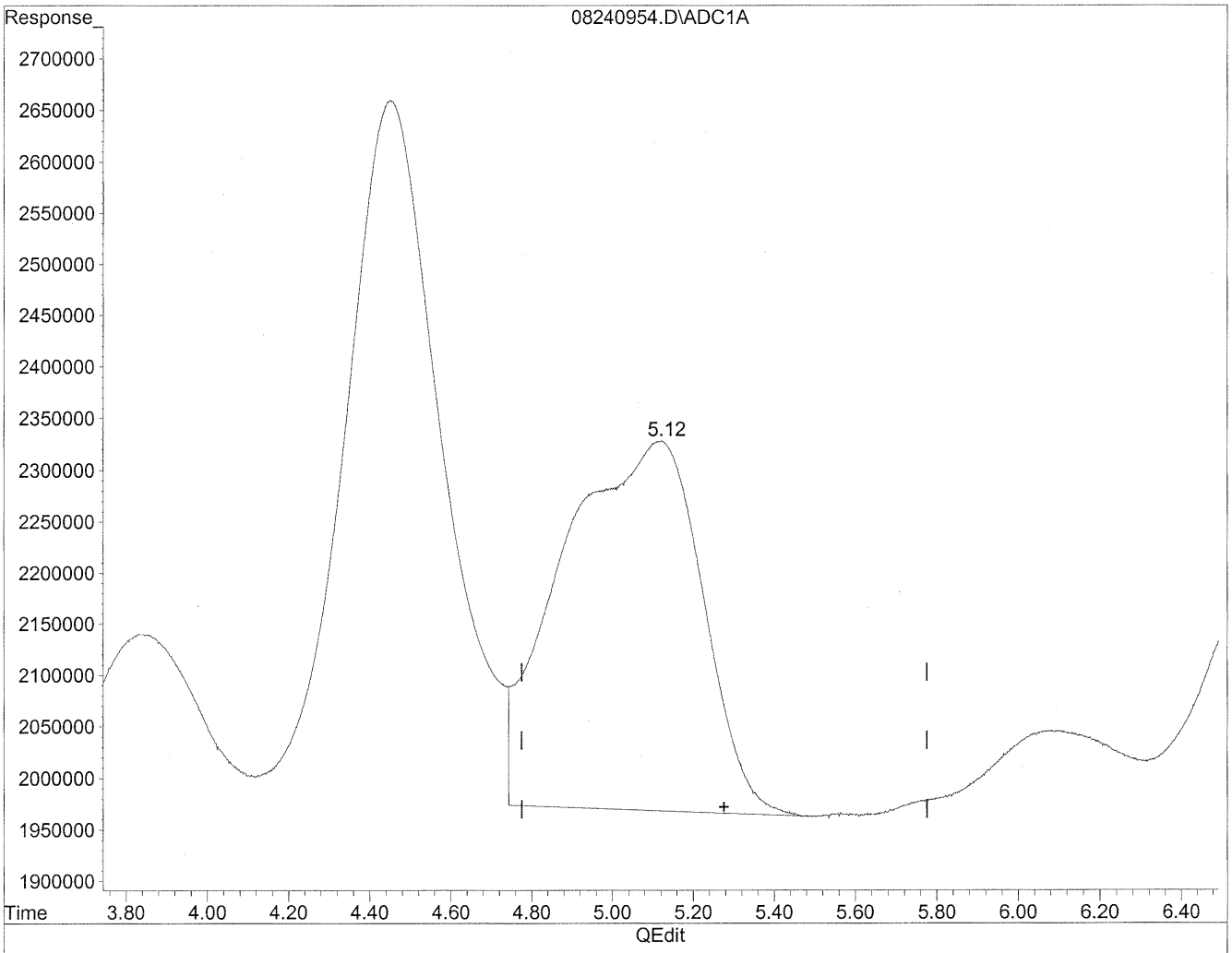
(+) = Expected Retention Time

08240954.D TO110709.M Sat Aug 29 14:45:00 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:56 19109 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



(5) Butyraldehyde

5.12min 978.608ng/ml

response 86446397

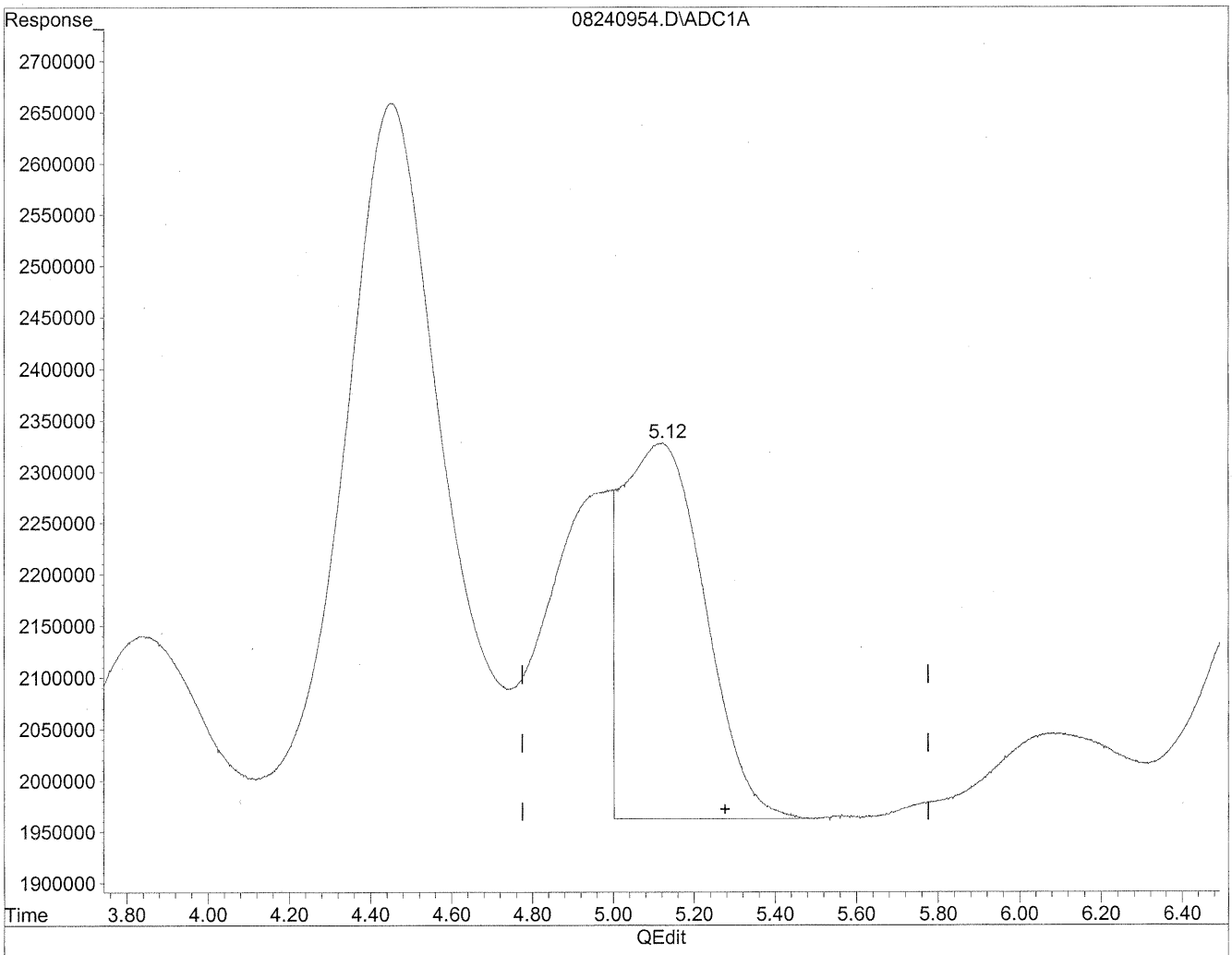
(+) = Expected Retention Time

08240954.D TO110709.M Sat Aug 29 14:45:04 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:56 19109 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



(5) Butyraldehyde

5.12min 587.491ng/ml m

response 51896644

HC 8/29/09 SP

15

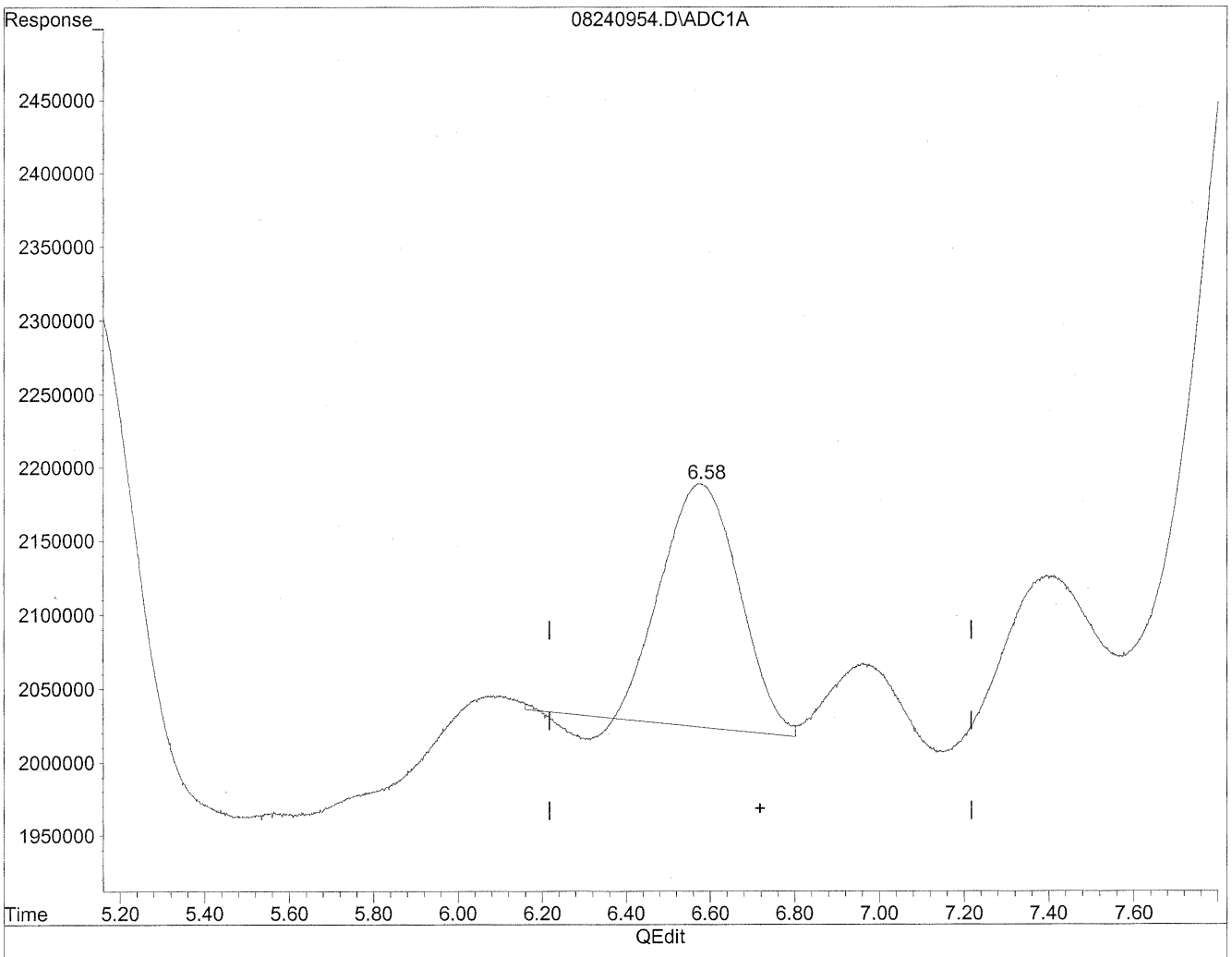
(+) = Expected Retention Time

08240954.D TO110709.M Sat Aug 29 14:45:28 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:56 19109 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



(6) Benzaldehyde

6.58min 311.406ng/ml

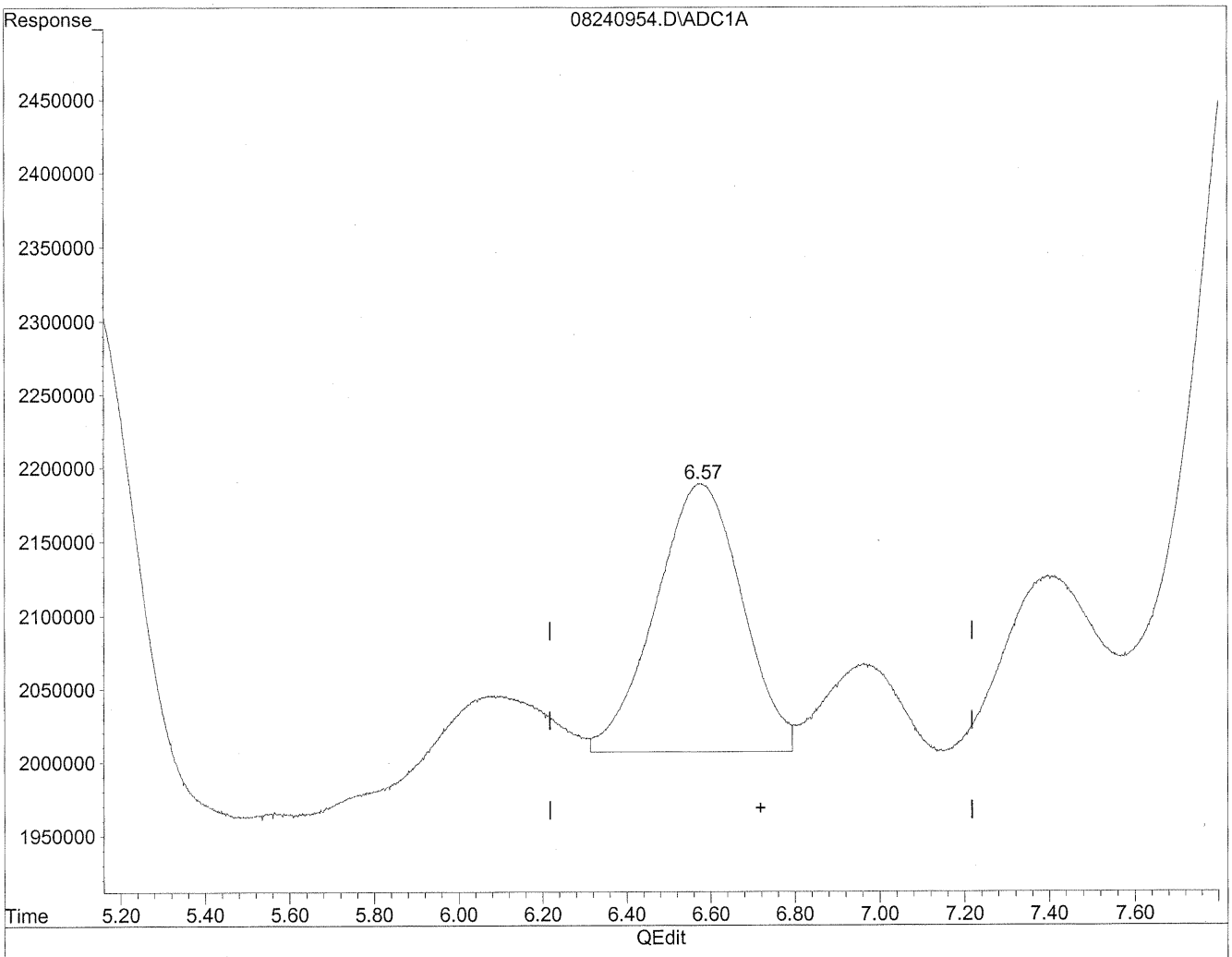
response 20512124

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:56 19109 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



(6) Benzaldehyde

6.57min 396.650ng/ml m

response 26127092

HC 8/31/09 RC

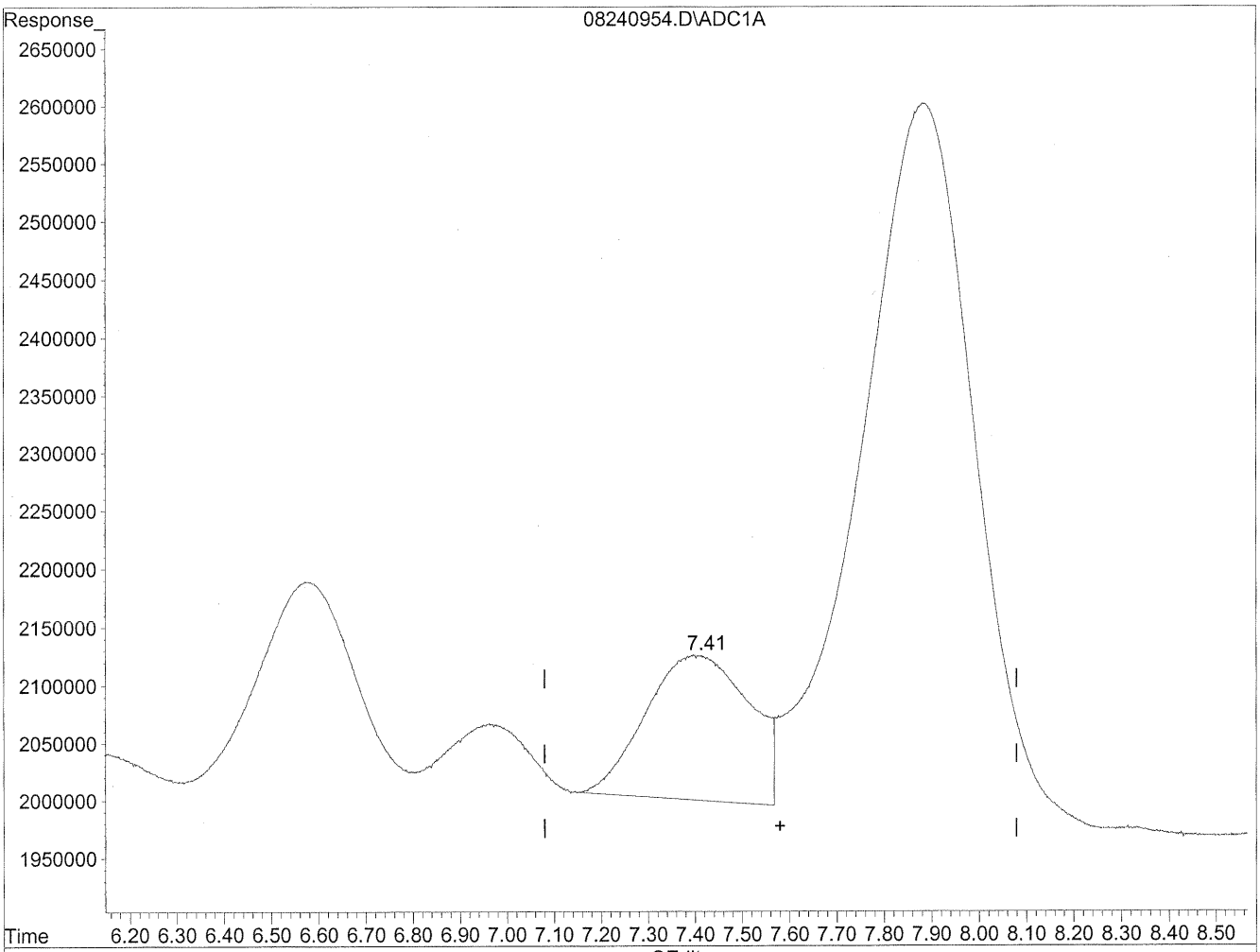
(+) = Expected Retention Time

08240954.D TO110709.M Sat Aug 29 14:46:22 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:56 19109 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



(7) Isovaleraldehyde

7.40min 246.774ng/ml

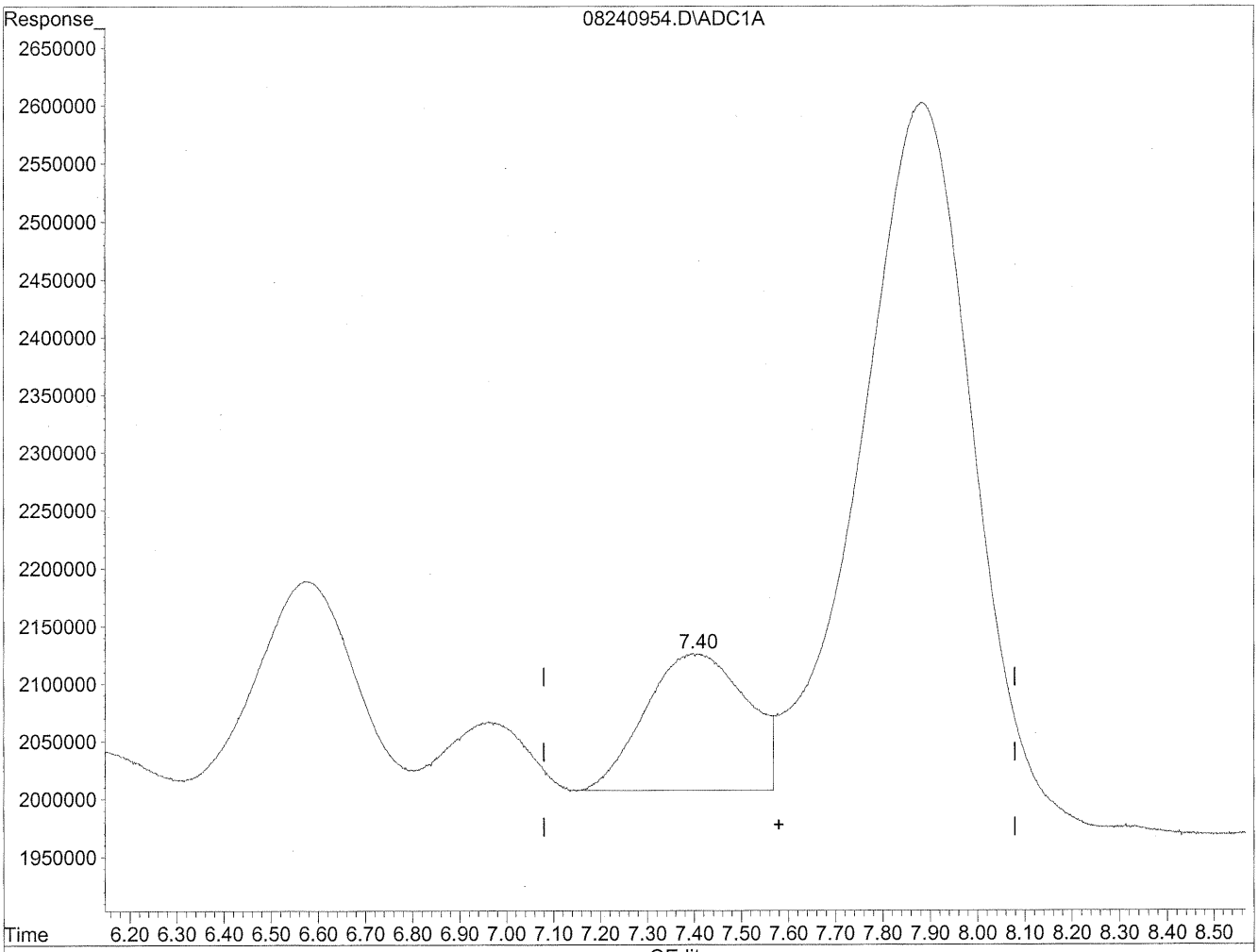
response 19310293

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:56 19109 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



(7) Isovaleraldehyde

7.40min 228.362ng/ml m

response 17869572

HC
8/29/09
BC

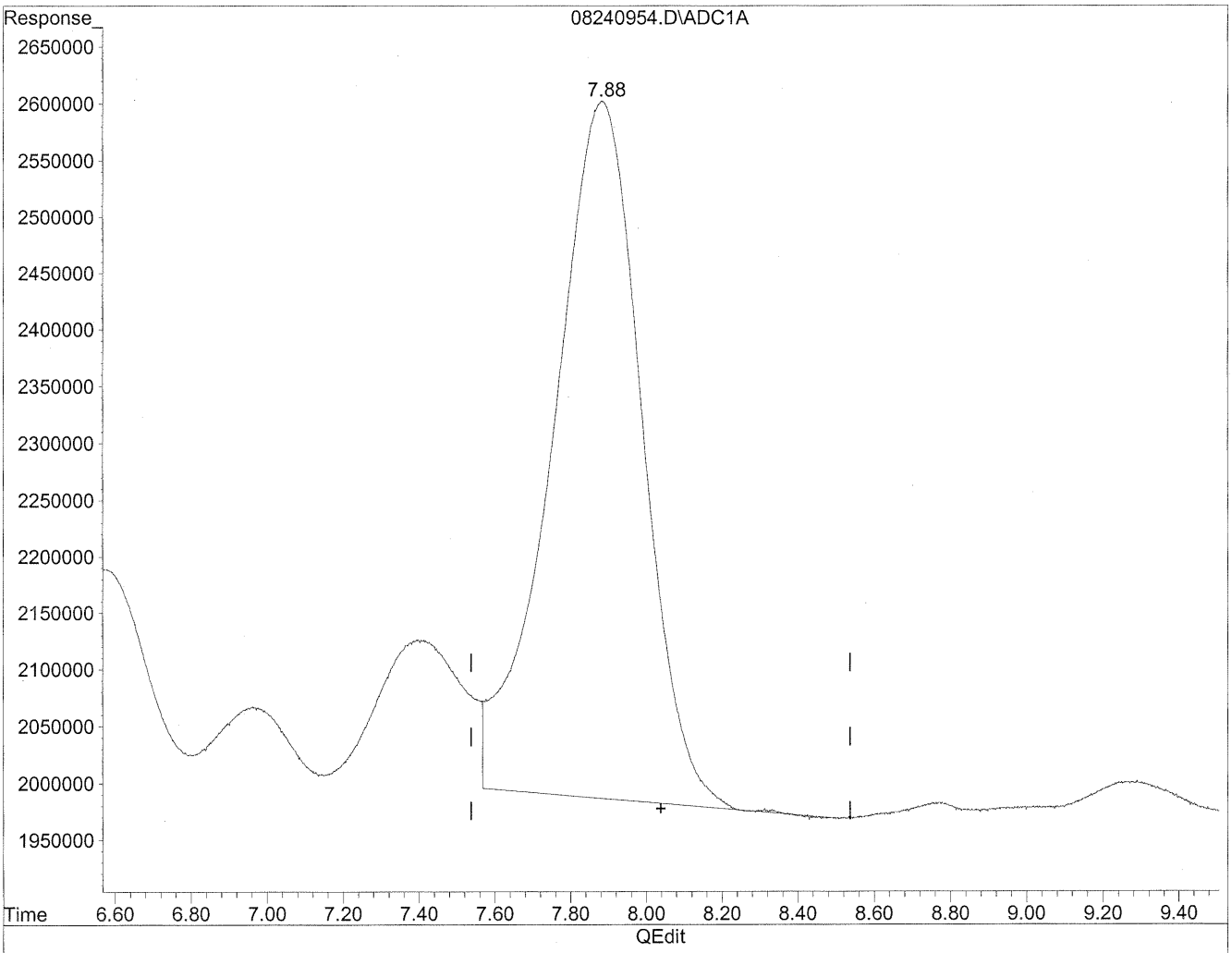
19

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:56 19109 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



(8) Valeraldehyde

7.88min 1376.868ng/ml

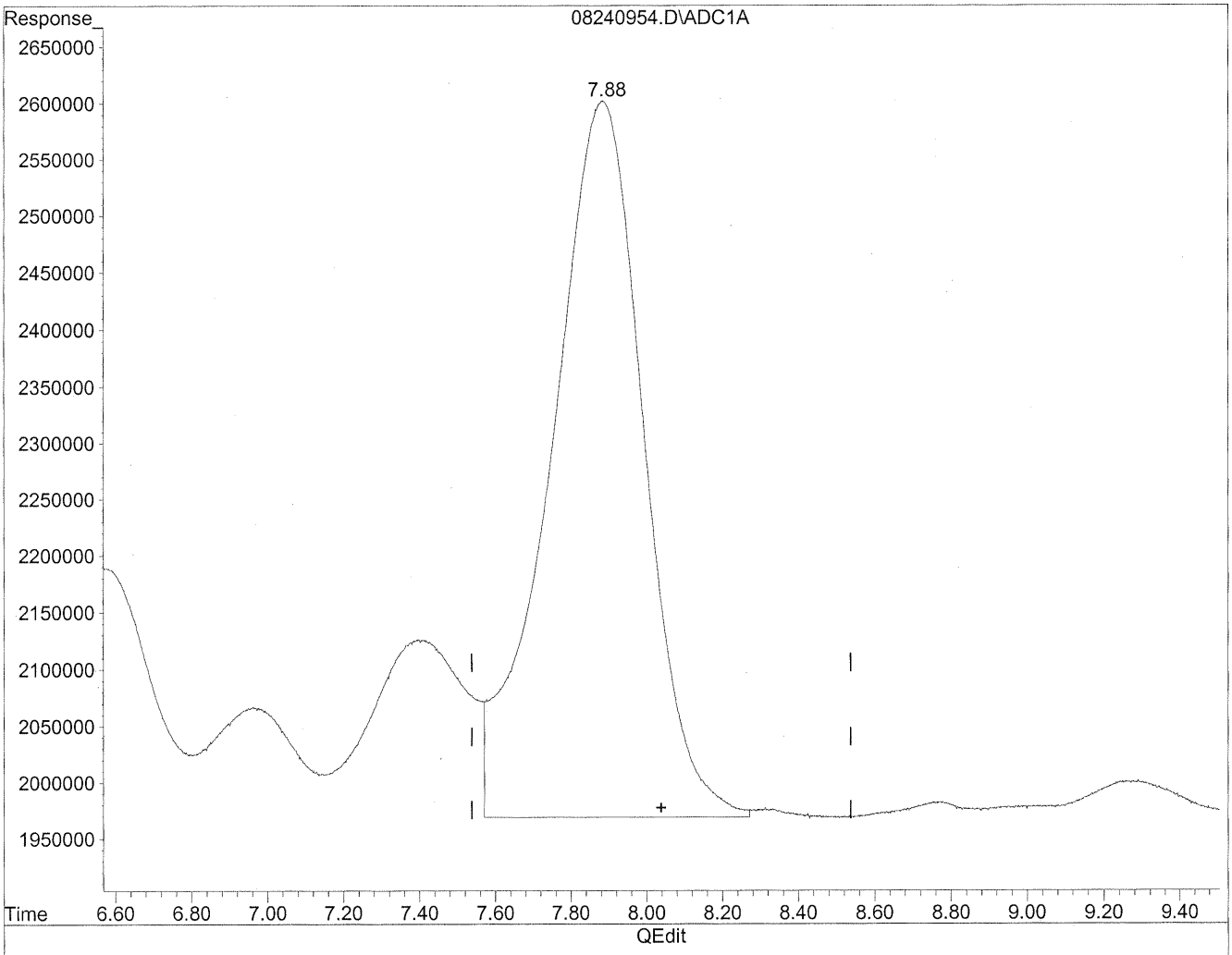
response 101206703

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240954.D Vial: 51
Acq On : 25 Aug 2009 1:47 am Operator: HC
Sample : P0902912-001 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:56 19109 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



(8) Valeraldehyde

7.88min 1468.104ng/ml m

response 107912963

HC 8/31/09 BC

21

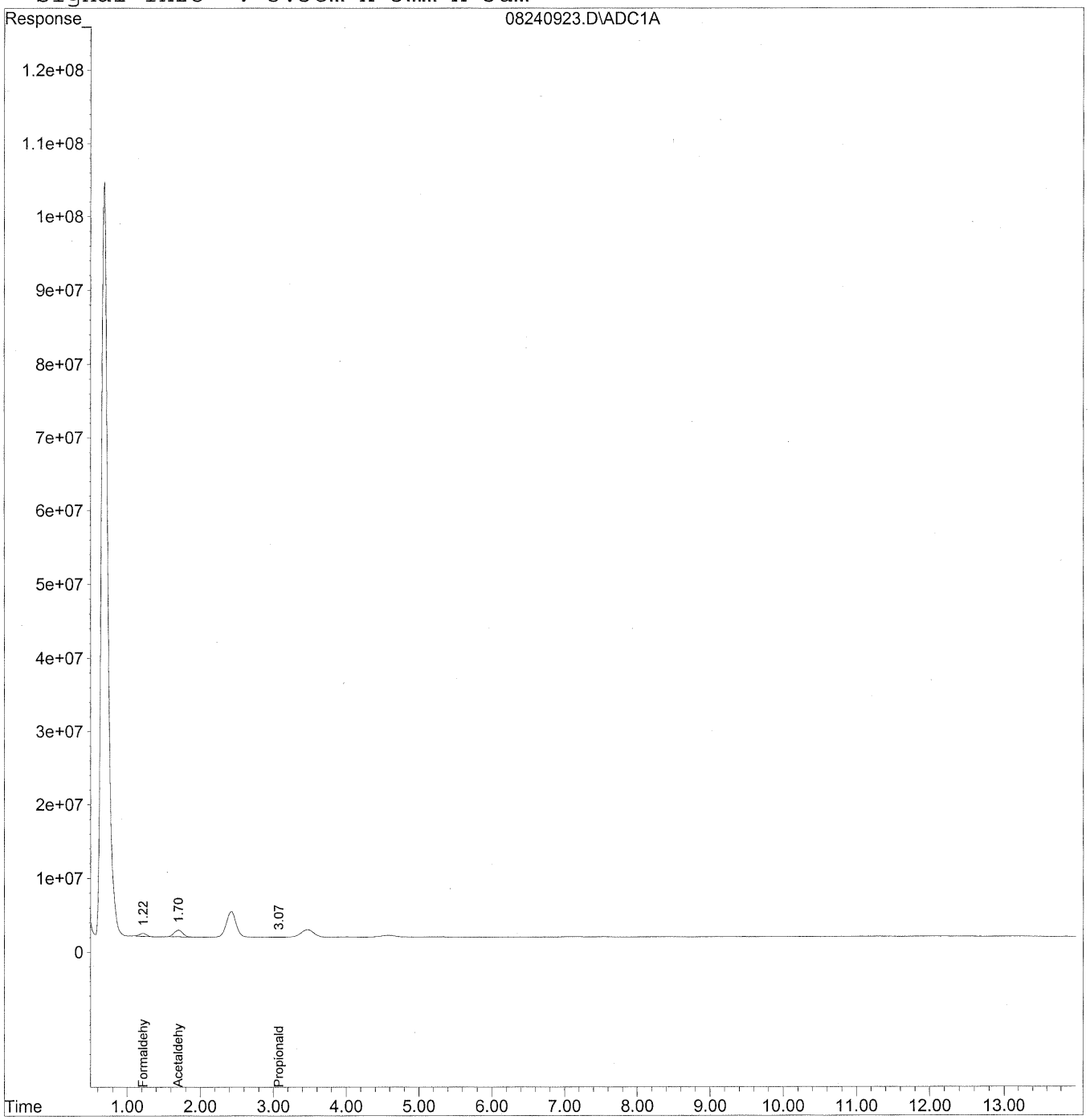
(+) = Expected Retention Time
08240954.D TO110709.M Sat Aug 29 14:47:52 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240923.D Vial: 21
Acq On : 24 Aug 2009 6:01 pm Operator: HC
Sample : P0902912-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 13:32 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 : Tue Aug 25 10:21:57 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

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



22

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\24\08240923.D Vial: 21
Acq On : 24 Aug 2009 6:01 pm Operator: HC
Sample : P0902912-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 13:32 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 : Tue Aug 25 10:21:57 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

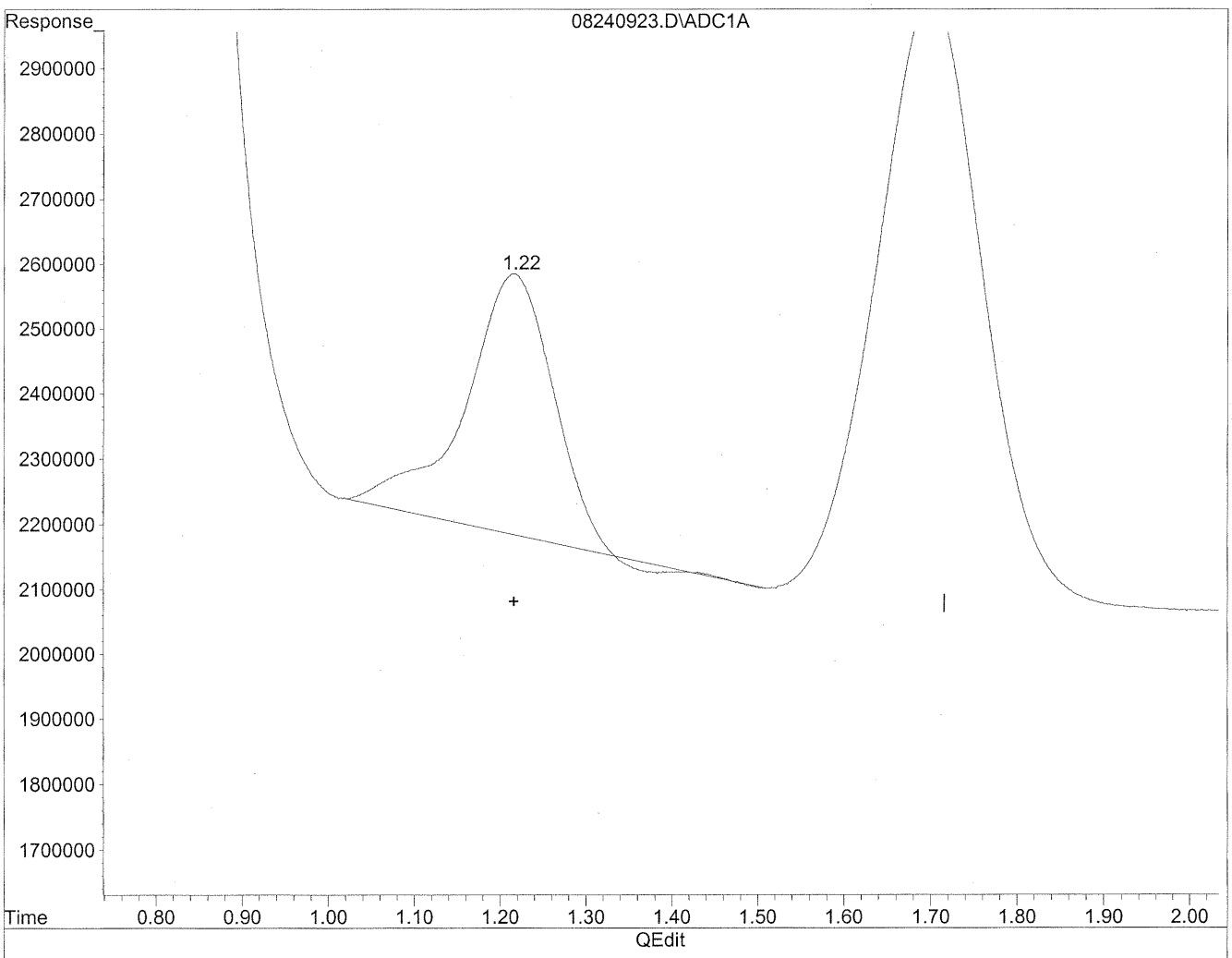
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	27267170	148.529	ng/mlm
2)	Acetaldehyde	1.70	78079421	556.821	ng/mlm
3)	Propionaldehyde	3.07	1950649	18.282	ng/mlm
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240923.D Vial: 21
Acq On : 24 Aug 2009 6:01 pm Operator: HC
Sample : P0902912-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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

1.22min 154.894ng/ml

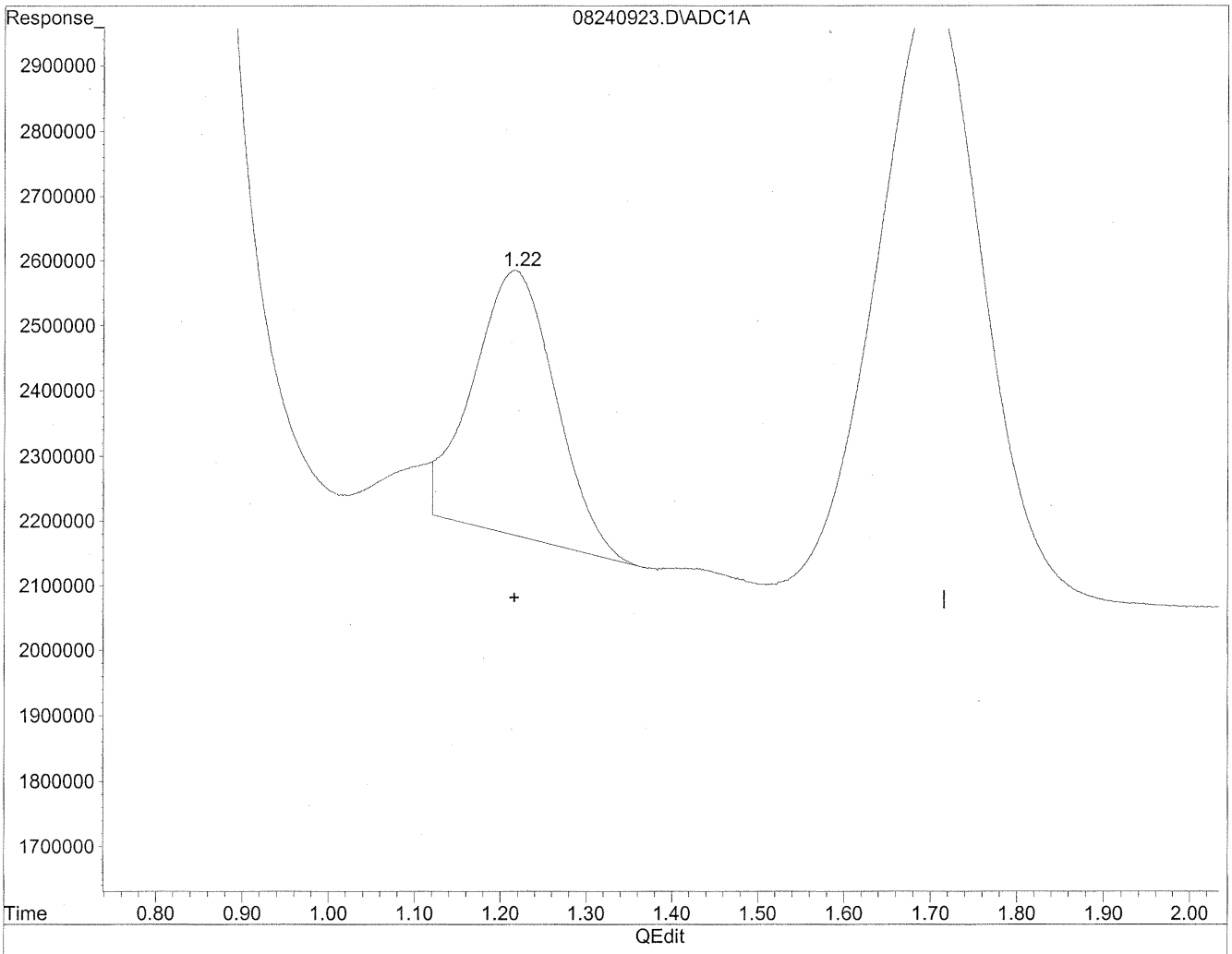
response 28435653

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240923.D Vial: 21
Acq On : 24 Aug 2009 6:01 pm Operator: HC
Sample : P0902912-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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

1.22min 148.529ng/ml m

response 27267170

HC 8/28/09 IC

8/31/09

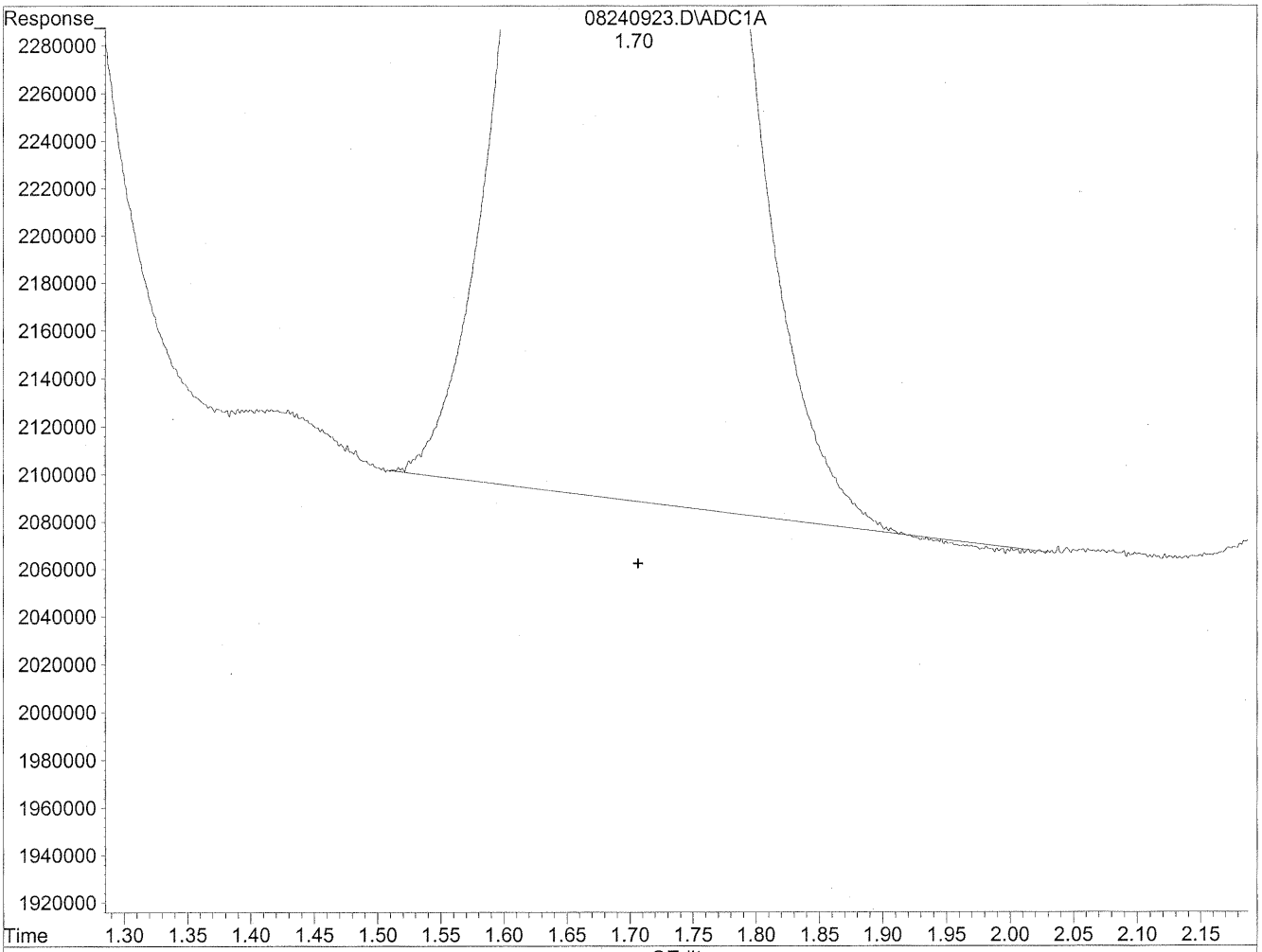
25

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240923.D Vial: 21
Acq On : 24 Aug 2009 6:01 pm Operator: HC
Sample : P0902912-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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 555.684ng/ml

response 77920011

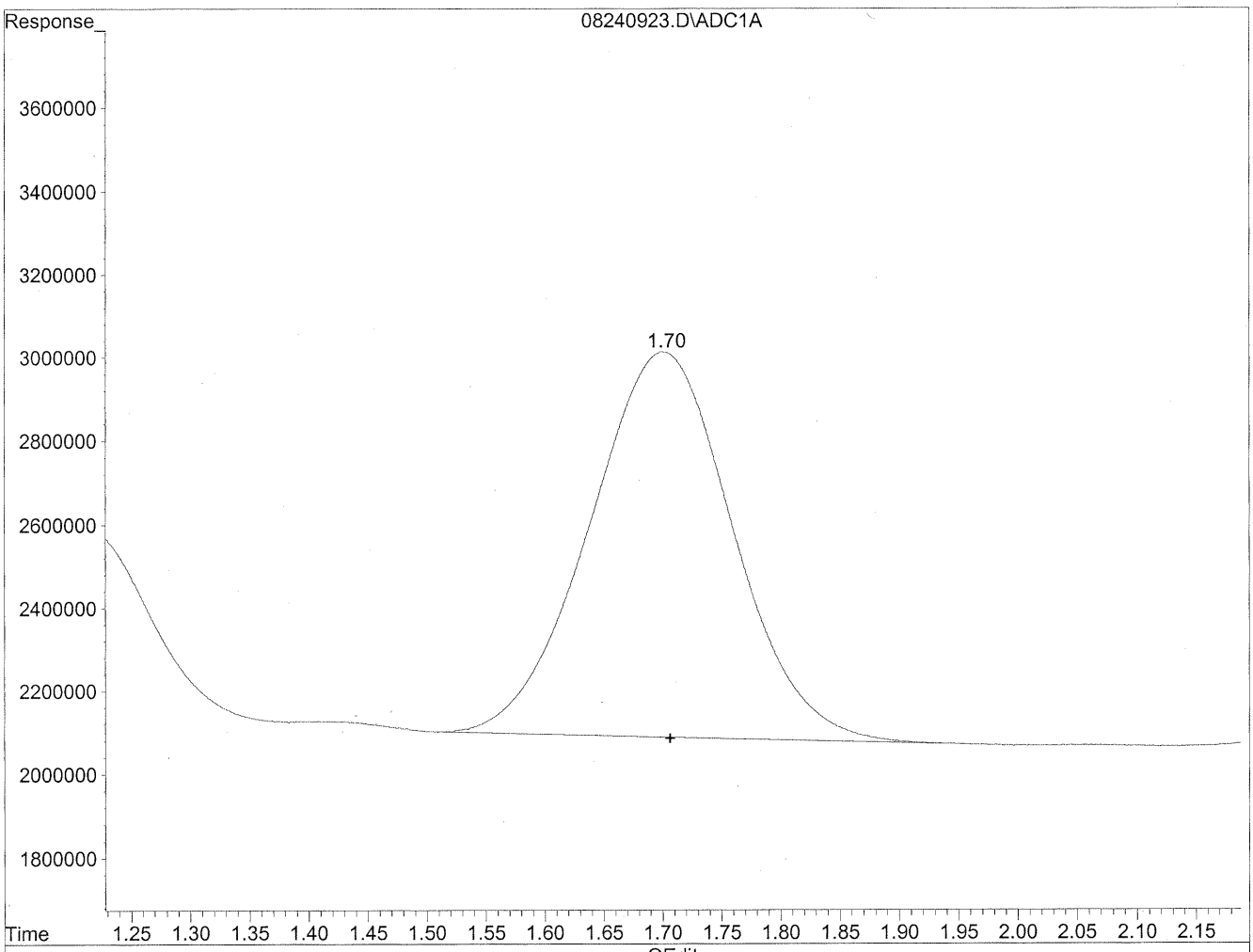
(+) = Expected Retention Time

08240923.D TO110709.M Sat Aug 29 13:31:36 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240923.D Vial: 21
Acq On : 24 Aug 2009 6:01 pm Operator: HC
Sample : P0902912-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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 556.821ng/ml m

response 78079421

HC 8/29/09 LC

8/29/09

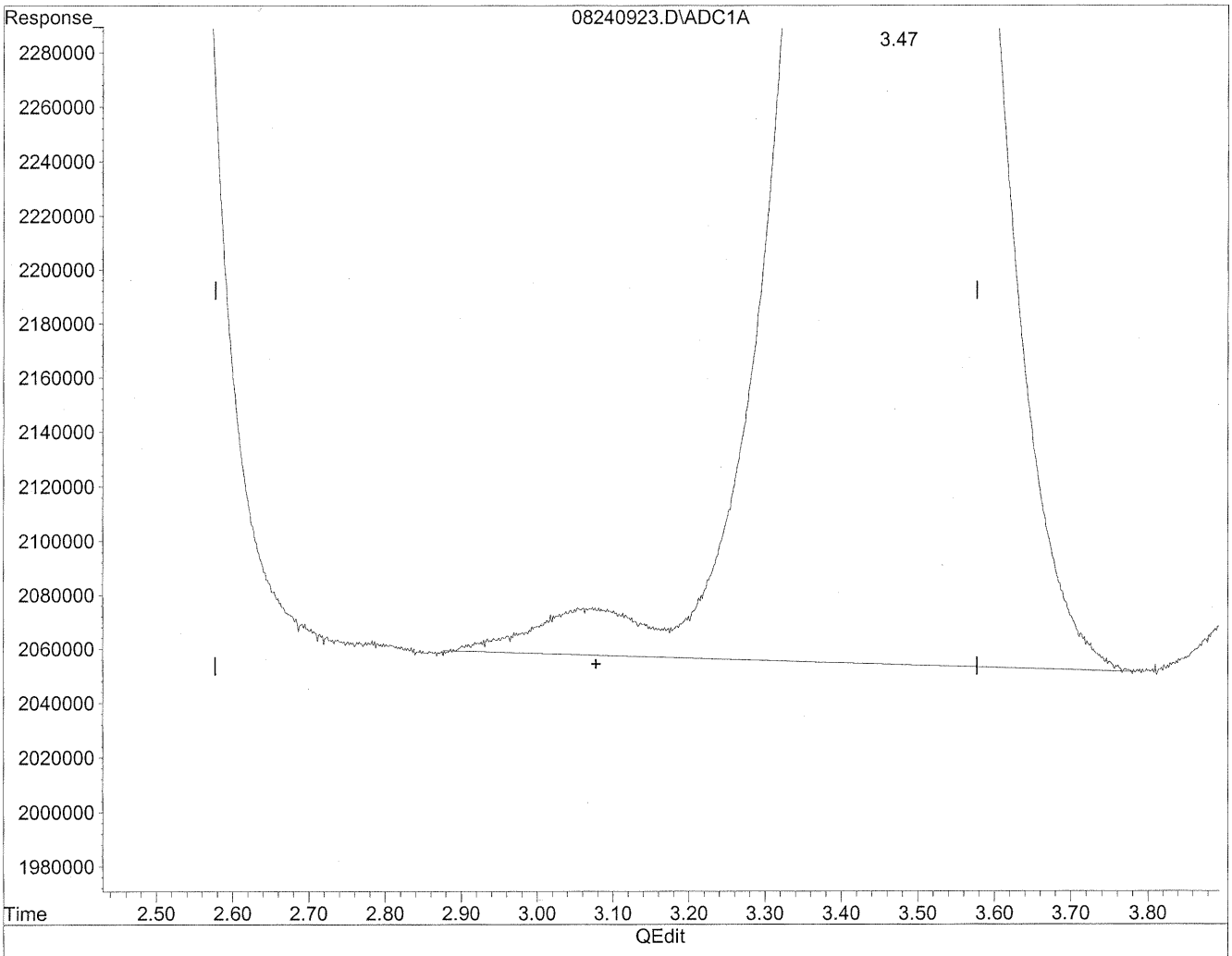
(+) = Expected Retention Time

08240923.D TO110709.M Sat Aug 29 13:31:46 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240923.D Vial: 21
Acq On : 24 Aug 2009 6:01 pm Operator: HC
Sample : P0902912-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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



(3) Propionaldehyde

3.47min 1209.870ng/ml

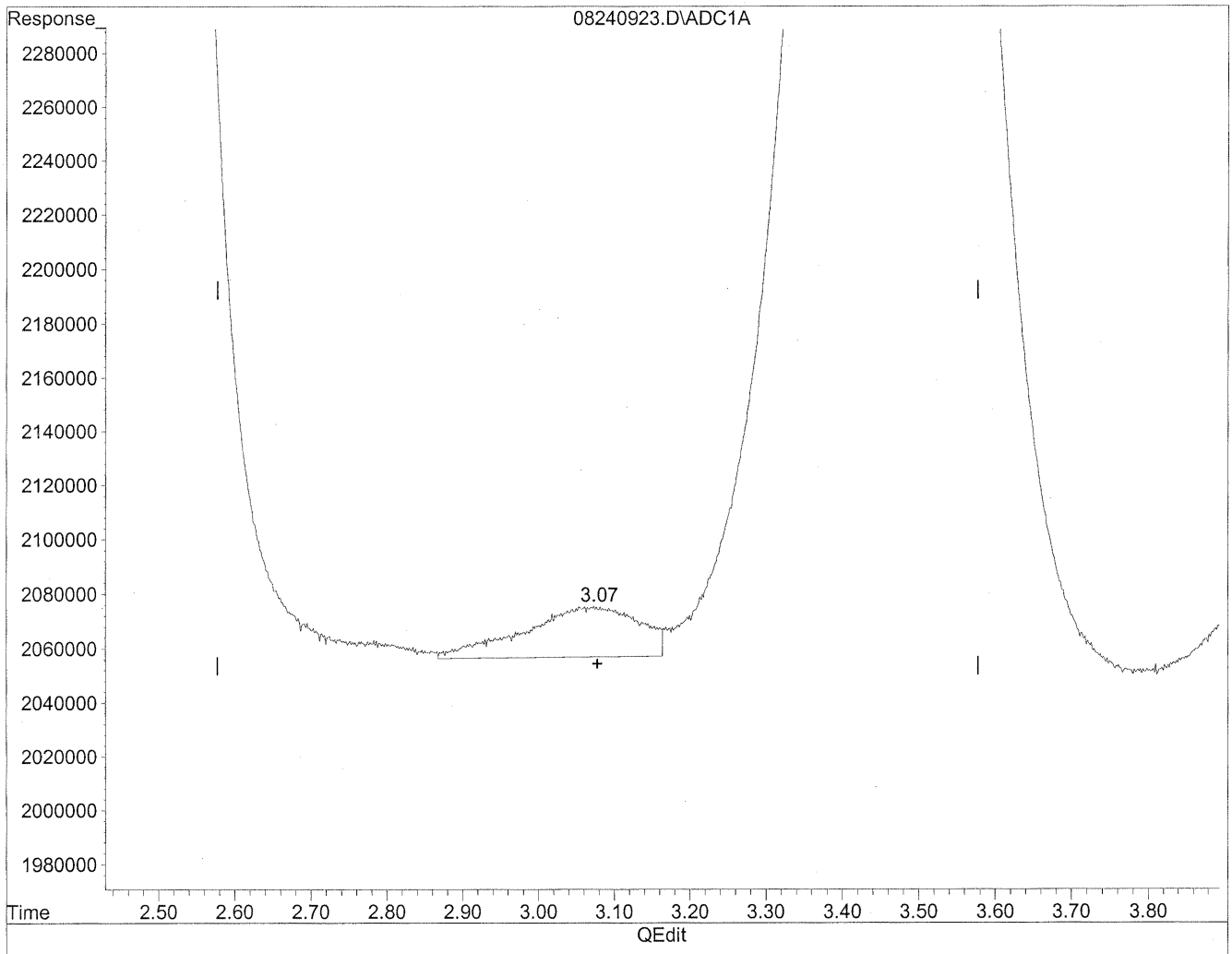
response 129087370

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240923.D Vial: 21
Acq On : 24 Aug 2009 6:01 pm Operator: HC
Sample : P0902912-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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



(3) Propionaldehyde

3.07min 18.282ng/ml m

response 1950649

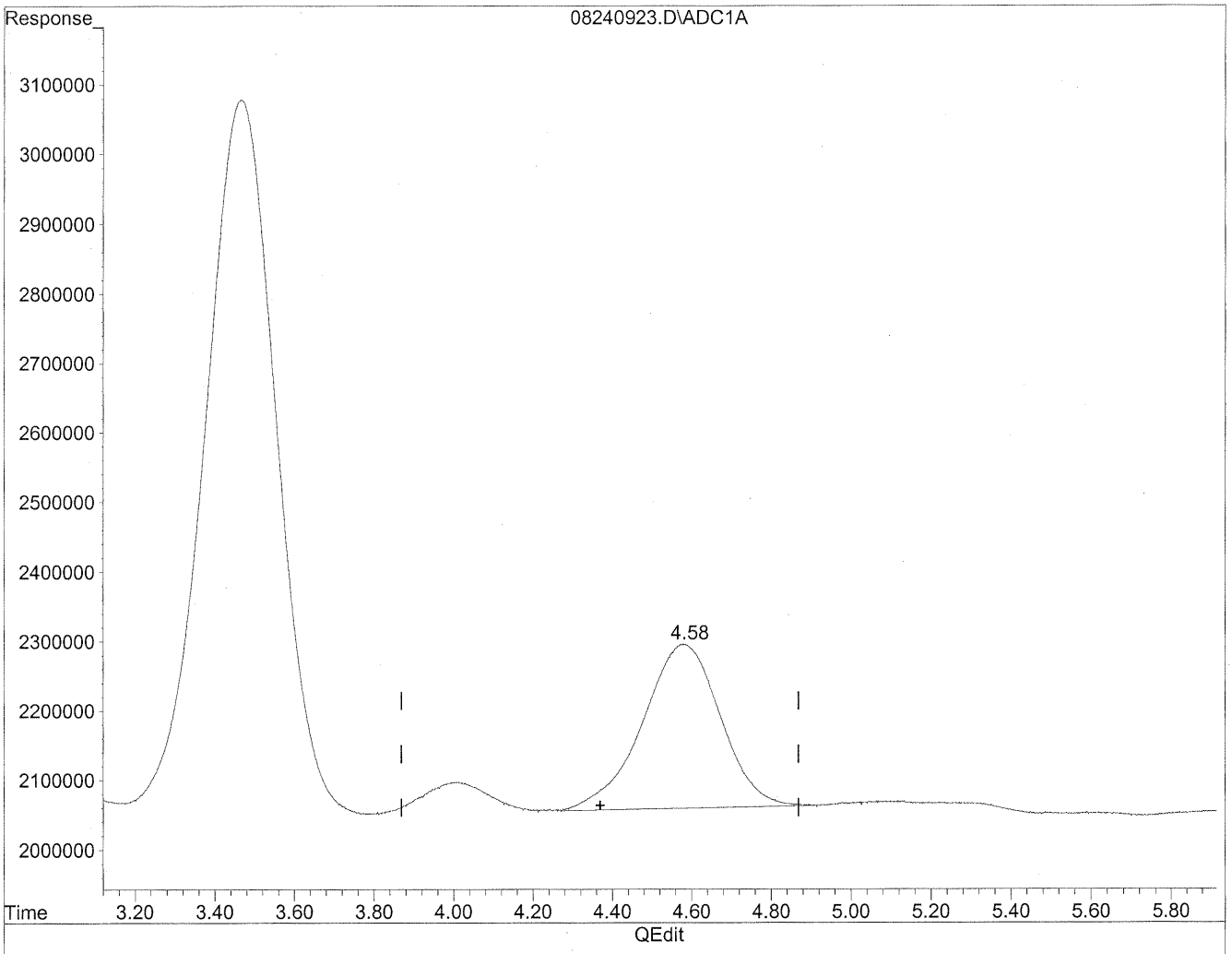
(+) = Expected Retention Time

08240923.D TO110709.M Sat Aug 29 13:31:59 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240923.D Vial: 21
Acq On : 24 Aug 2009 6:01 pm Operator: HC
Sample : P0902912-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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



(4) Crotonaldehyde

4.58min 332.167ng/ml

response 32358096

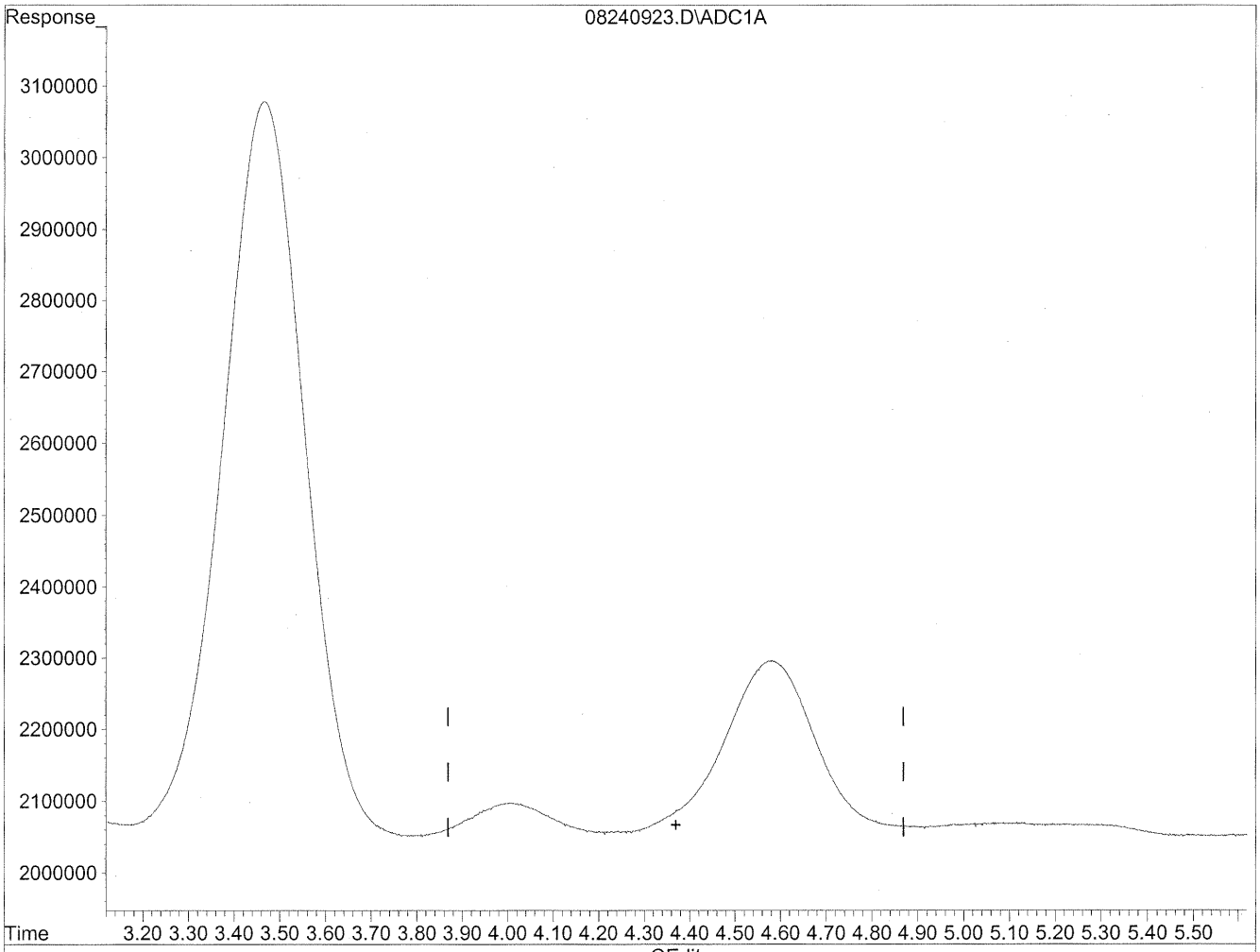
(+) = Expected Retention Time

08240923.D TO110709.M Sat Aug 29 13:32:03 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240923.D Vial: 21
Acq On : 24 Aug 2009 6:01 pm Operator: HC
Sample : P0902912-001 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/31/09 WP 8/31/09

(+) = Expected Retention Time

08240923.D TO110709.M Sat Aug 29 13:32:07 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **102511**
Client Project ID: **16512**

CAS Project ID: P0902912
CAS Sample ID: P0902912-002

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: 8/20/09
Date Received: 8/21/09
Date Analyzed: 8/25/09
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**	**330**	**3.2**	0.99	**2.6**	0.80	
75-07-0	Acetaldehyde	< 100	ND	0.99	ND	0.55	
123-38-6	Propionaldehyde	< 100	ND	0.99	ND	0.41	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.99	ND	0.34	
123-72-8	Butyraldehyde	< 100	ND	0.99	ND	0.33	
100-52-7	Benzaldehyde	< 100	ND	0.99	ND	0.23	
590-86-3	Isovaleraldehyde	< 100	ND	0.99	ND	0.28	
110-62-3	Valeraldehyde	< 100	ND	0.99	ND	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	< 100	ND	0.99	ND	0.24	
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	ND	0.99	ND	0.18	

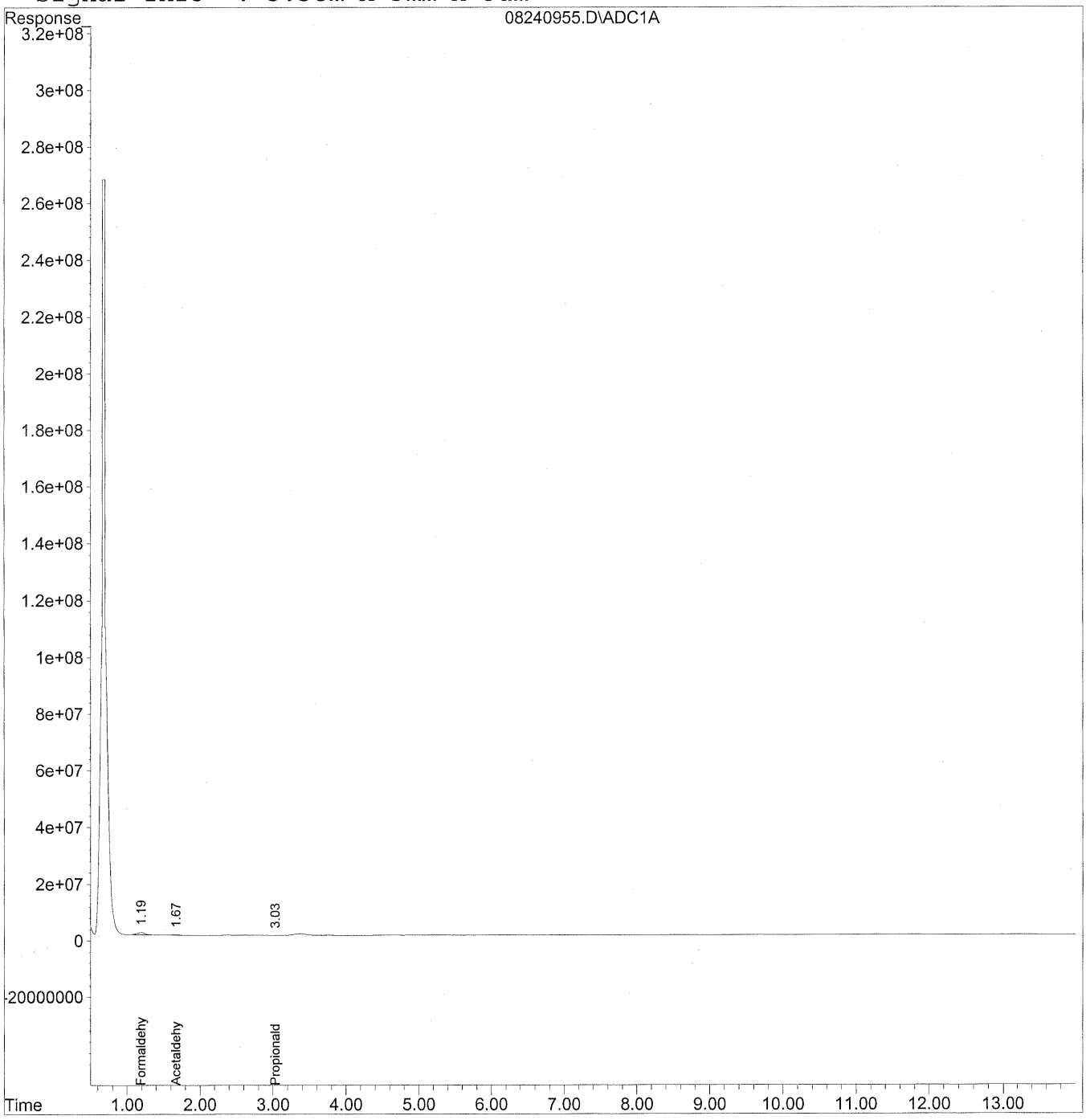
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.

Verified By: Ro Date: 9/2/09

Data File : J:\LC01\DATA\TO11\2009_08\24\08240955.D Vial: 52
Acq On : 25 Aug 2009 2:02 am Operator: HC
Sample : P0902912-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 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 : Mon Aug 24 08:44:34 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\24\08240955.D Vial: 52
Acq On : 25 Aug 2009 2:02 am Operator: HC
Sample : P0902912-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 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 : Mon Aug 24 08:44:34 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

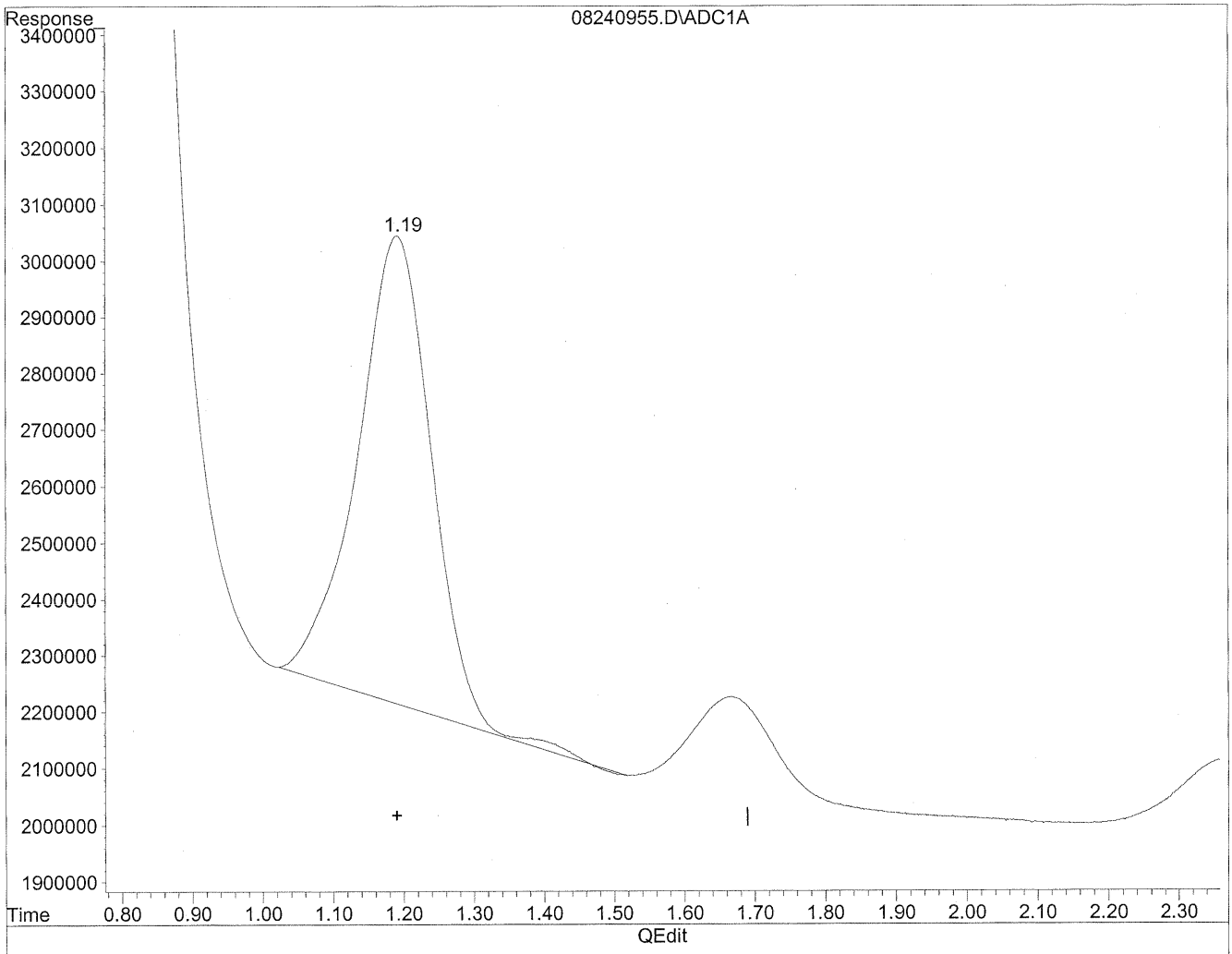
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.19	60333353	328.646	ng/mlm
2)	Acetaldehyde	1.67	12813955	91.382	ng/mlm
3)	Propionaldehyde	3.03	1078954	10.112	ng/mlm
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/mld
7)	Isovaleraldehyde	0.00	0	N.D.	ng/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\24\08240955.D Vial: 52
Acq On : 25 Aug 2009 2:02 am Operator: HC
Sample : P0902912-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.19min 333.119ng/ml

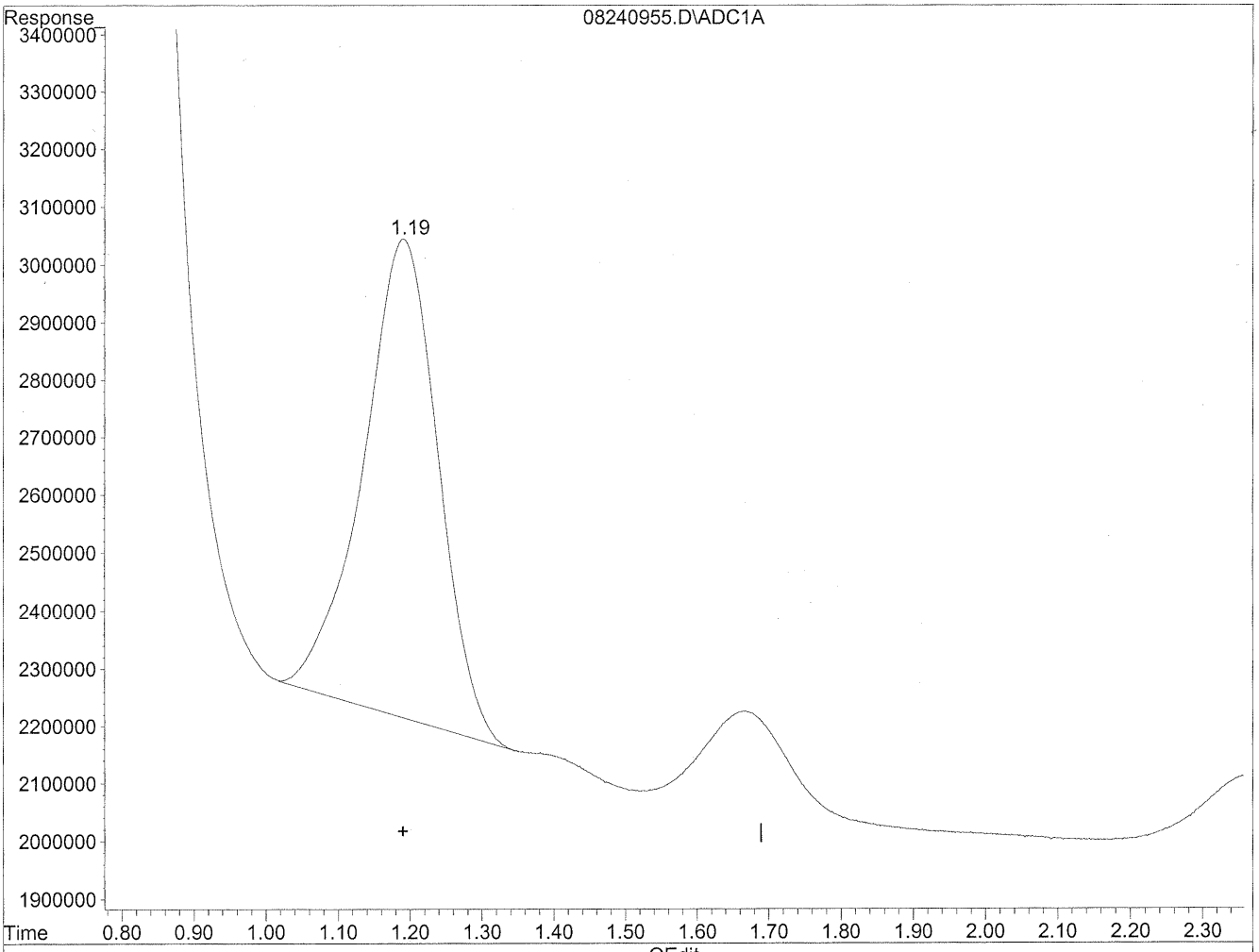
response 61154367

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240955.D Vial: 52
Acq On : 25 Aug 2009 2:02 am Operator: HC
Sample : P0902912-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.19min 328.646ng/ml m

response 60333353

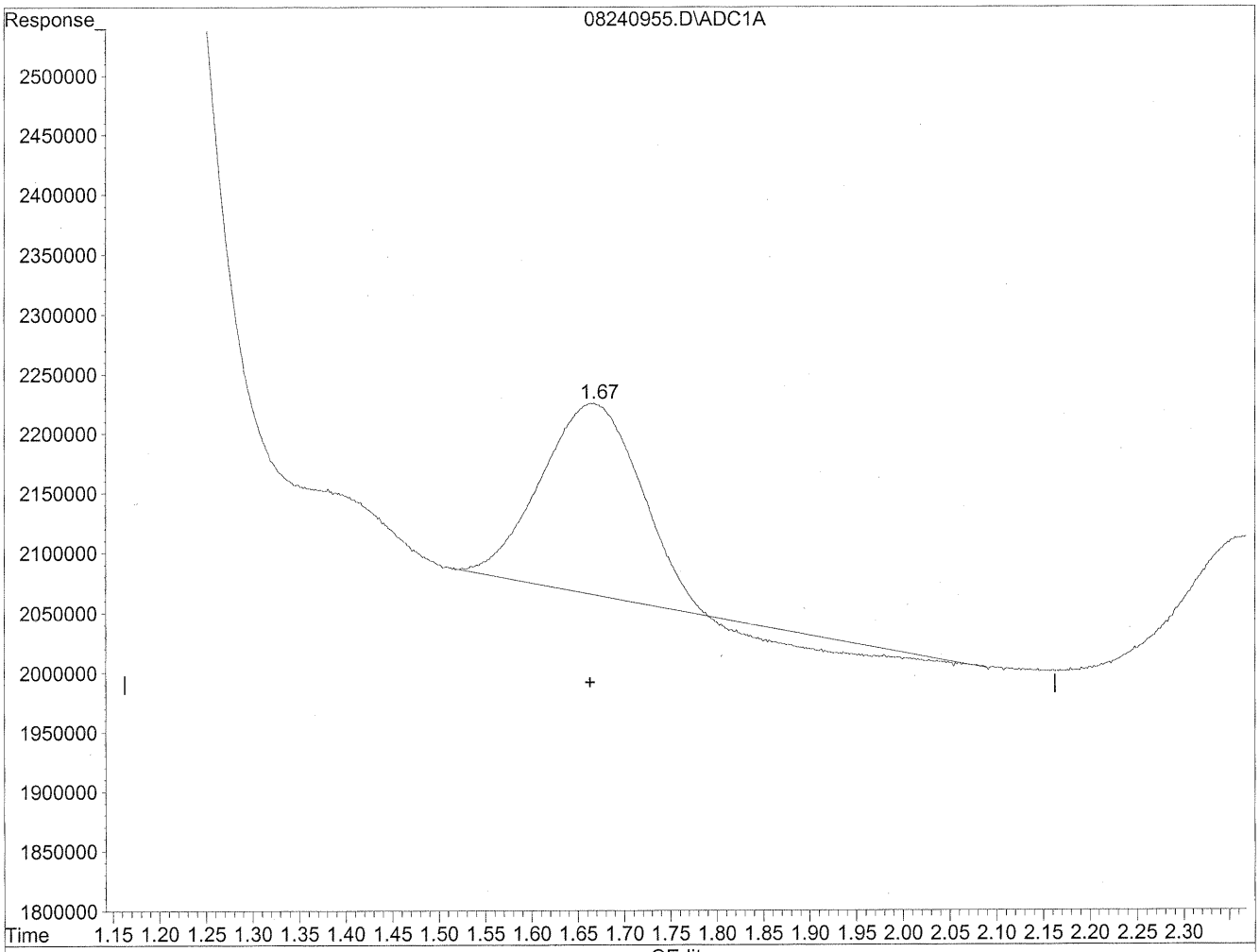
HC
8/26/09
IC

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240955.D Vial: 52
Acq On : 25 Aug 2009 2:02 am Operator: HC
Sample : P0902912-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.67min 77.636ng/ml

response 10886341

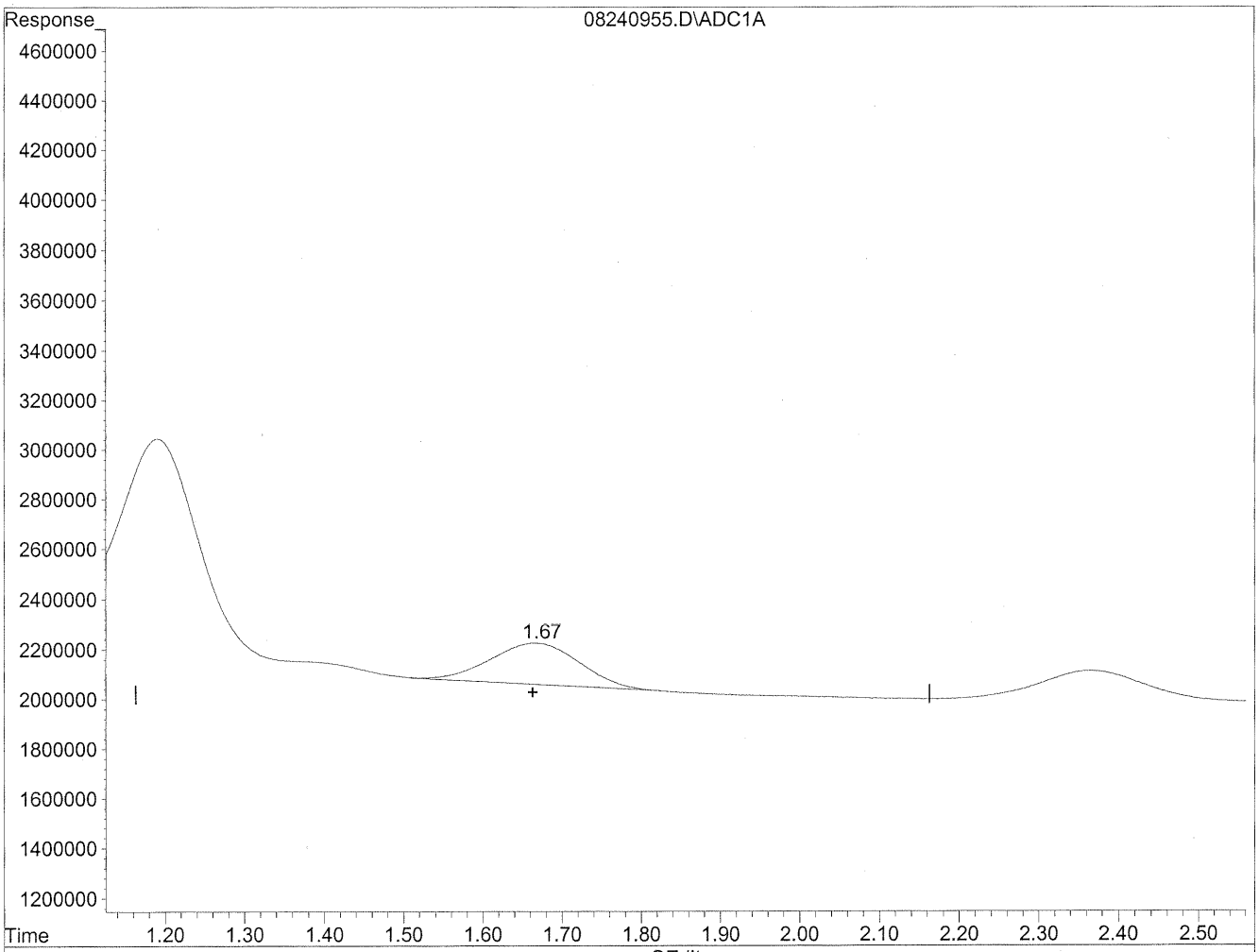
(+) = Expected Retention Time

08240955.D TO110709.M Sat Aug 29 14:48:33 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240955.D Vial: 52
Acq On : 25 Aug 2009 2:02 am Operator: HC
Sample : P0902912-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



QEdit

(2) Acetaldehyde

1.67min 91.382ng/ml m

response 12813955

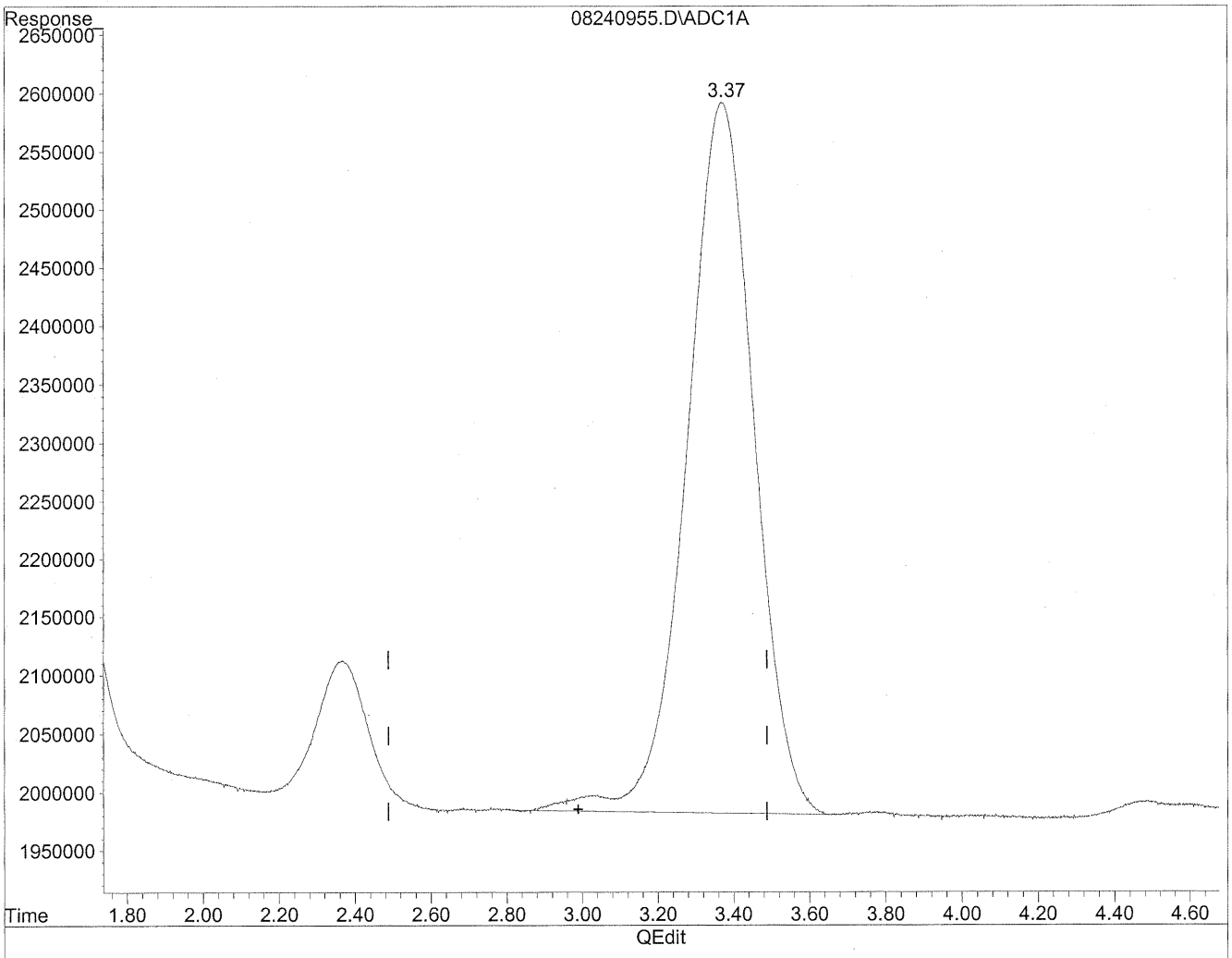
(+) = Expected Retention Time

08240955.D TO110709.M Sat Aug 29 14:48:40 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240955.D Vial: 52
Acq On : 25 Aug 2009 2:02 am Operator: HC
Sample : P0902912-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.37min 703.428ng/ml

response 75052384

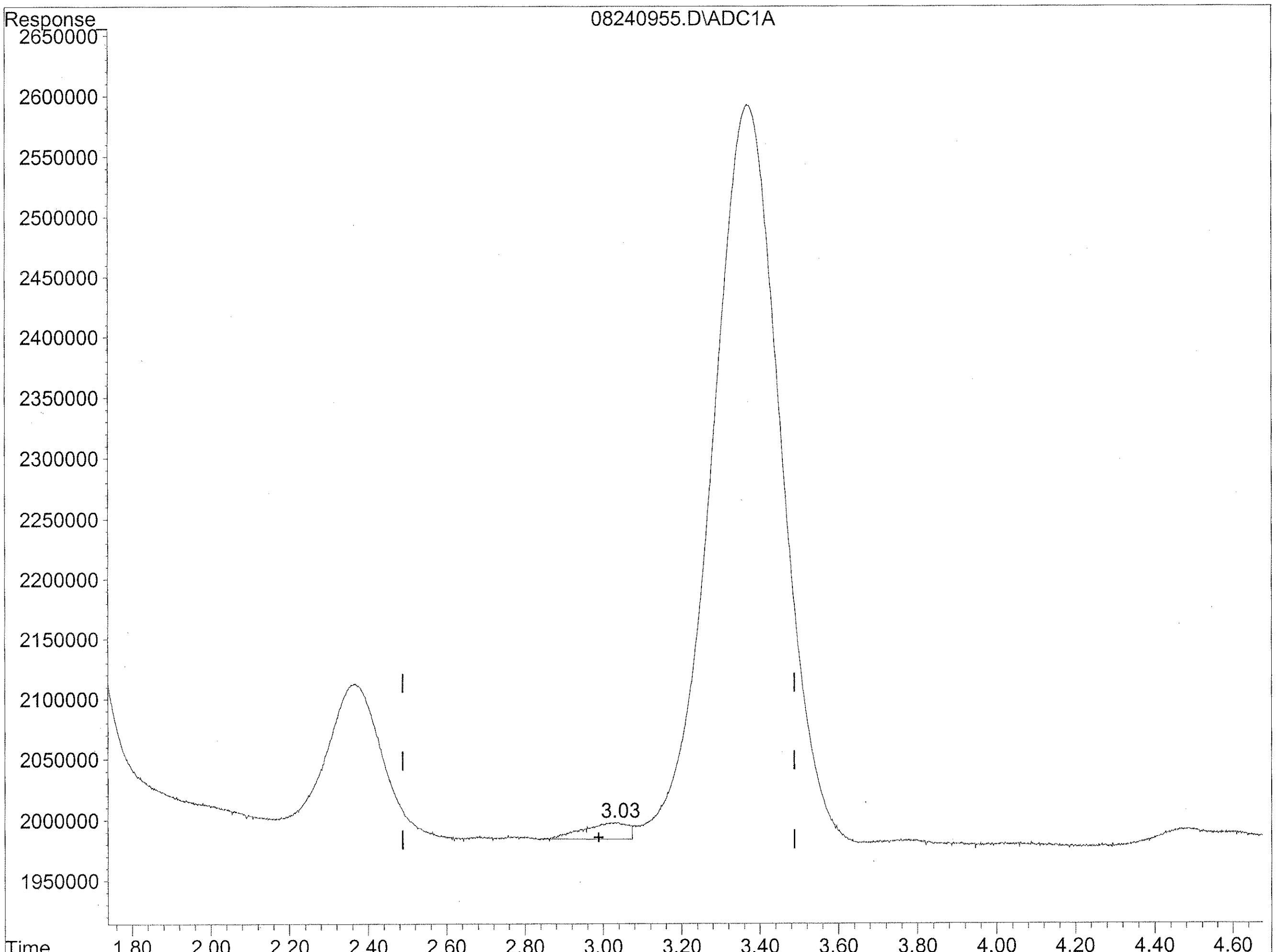
(+) = Expected Retention Time

08240955.D TO110709.M Sat Aug 29 14:48:48 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240955.D Vial: 52
Acq On : 25 Aug 2009 2:02 am Operator: HC
Sample : P0902912-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



QEdit

(3) Propionaldehyde

3.03min 10.112ng/ml m

response 1078954

HC 8/26/09

WP

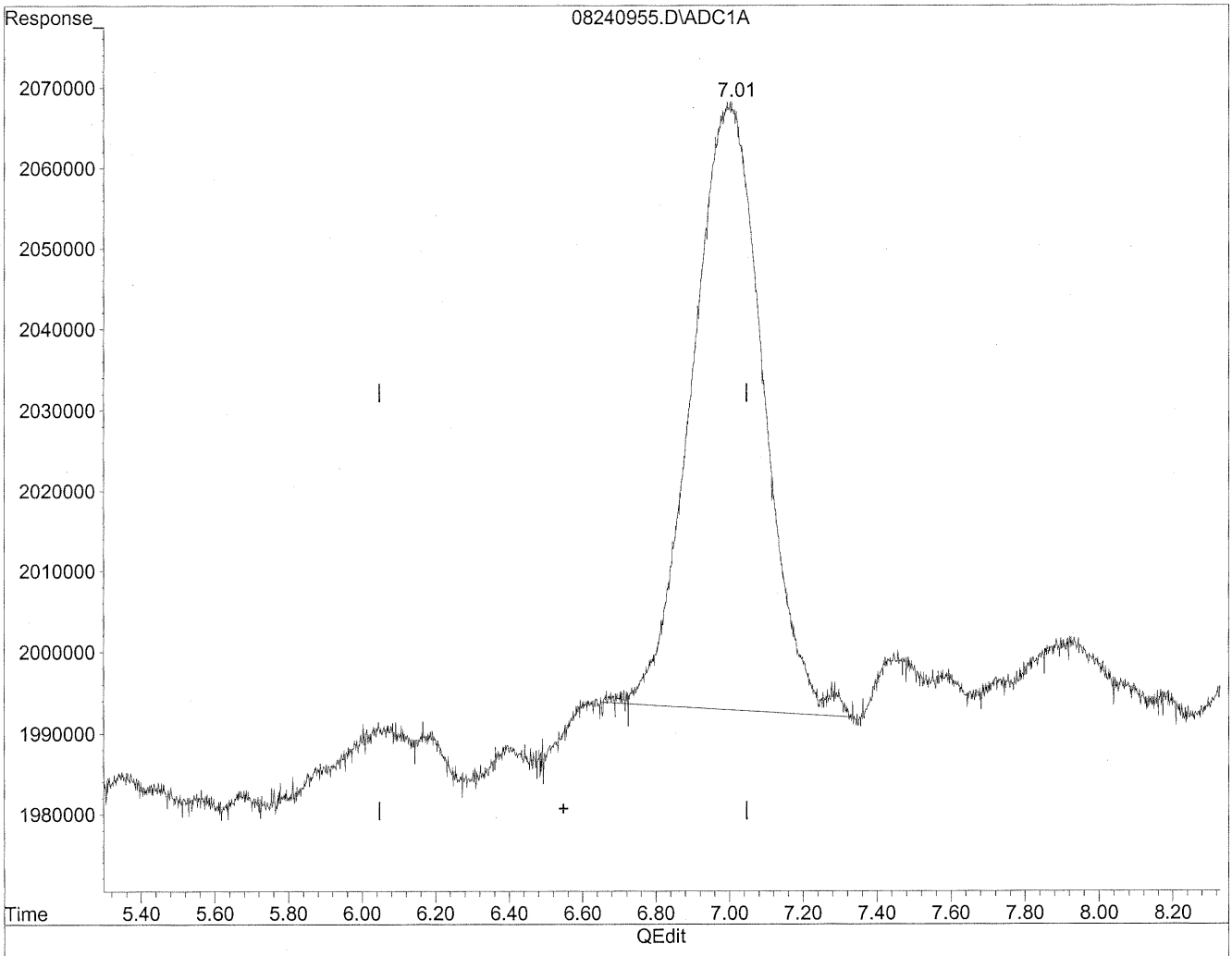
VH 8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240955.D Vial: 52
Acq On : 25 Aug 2009 2:02 am Operator: HC
Sample : P0902912-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

7.00min 153.064ng/ml

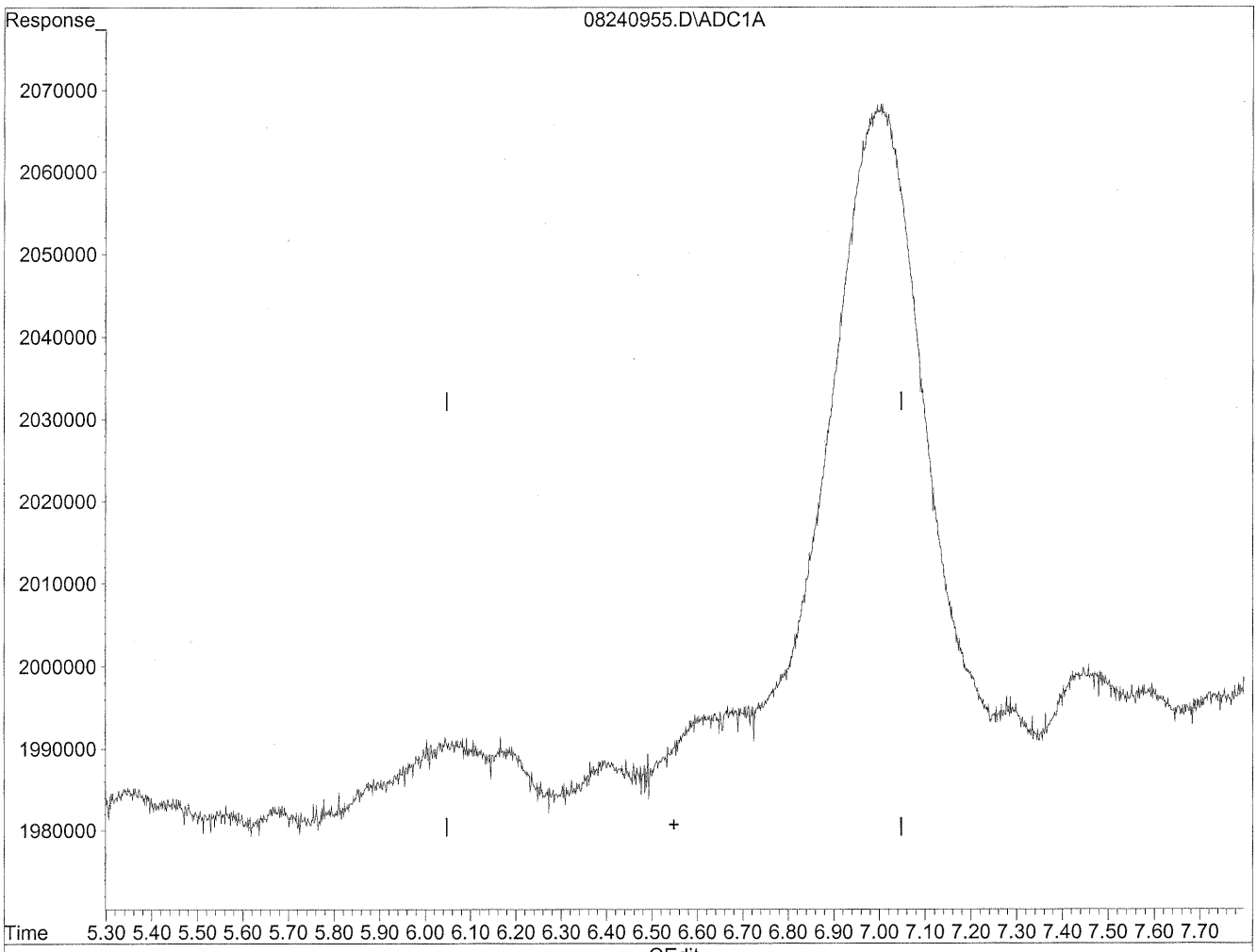
response 10082246

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240955.D Vial: 52
Acq On : 25 Aug 2009 2:02 am Operator: HC
Sample : P0902912-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/31/09

8/31/09

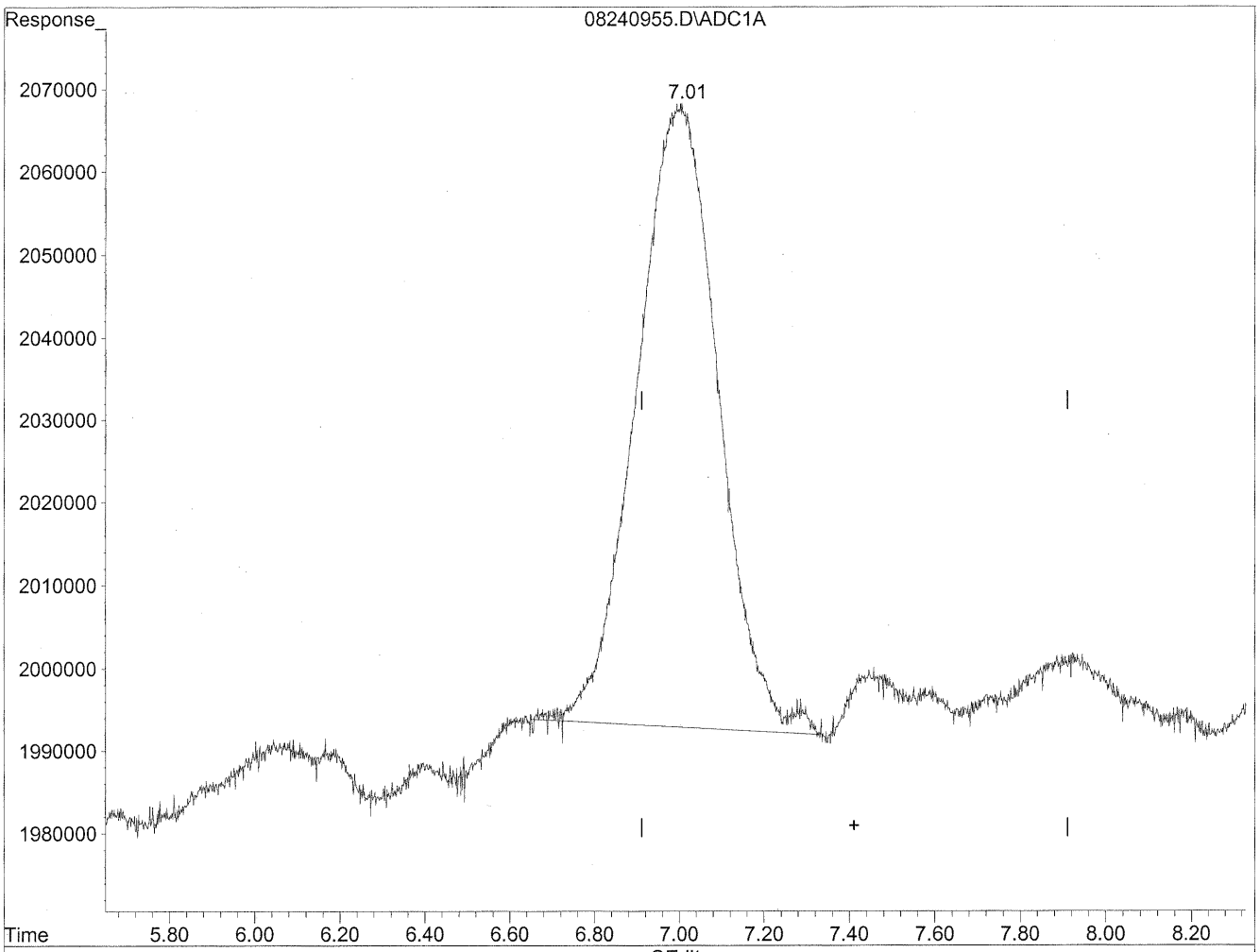
(+) = Expected Retention Time

08240955.D TO110709.M Sat Aug 29 14:49:06 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240955.D Vial: 52
Acq On : 25 Aug 2009 2:02 am Operator: HC
Sample : P0902912-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.00min 128.845ng/ml

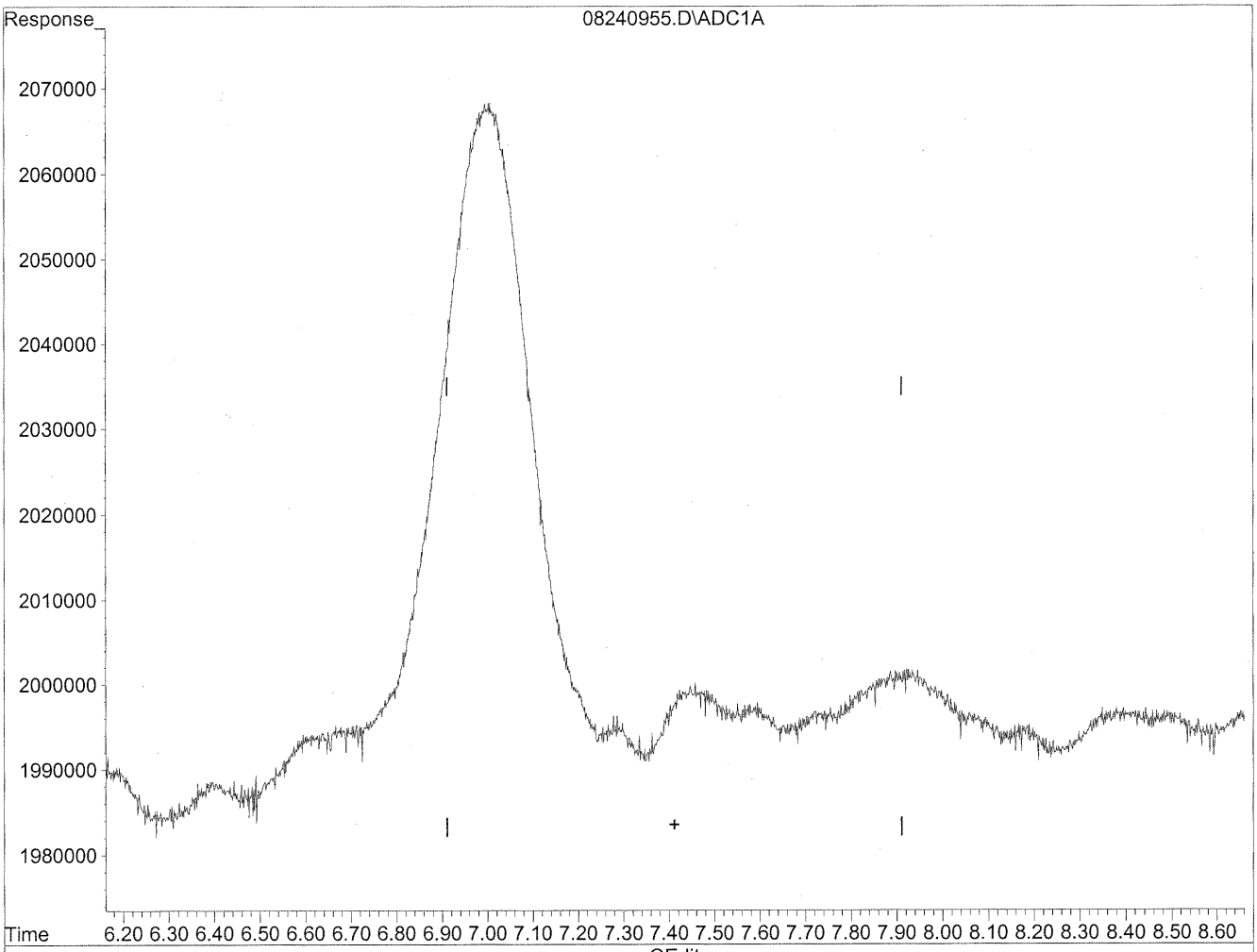
response 10082246

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240955.D Vial: 52
Acq On : 25 Aug 2009 2:02 am Operator: HC
Sample : P0902912-002 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

0.00min 0.000ng/ml d

response 0

HC 8/29/09 WP

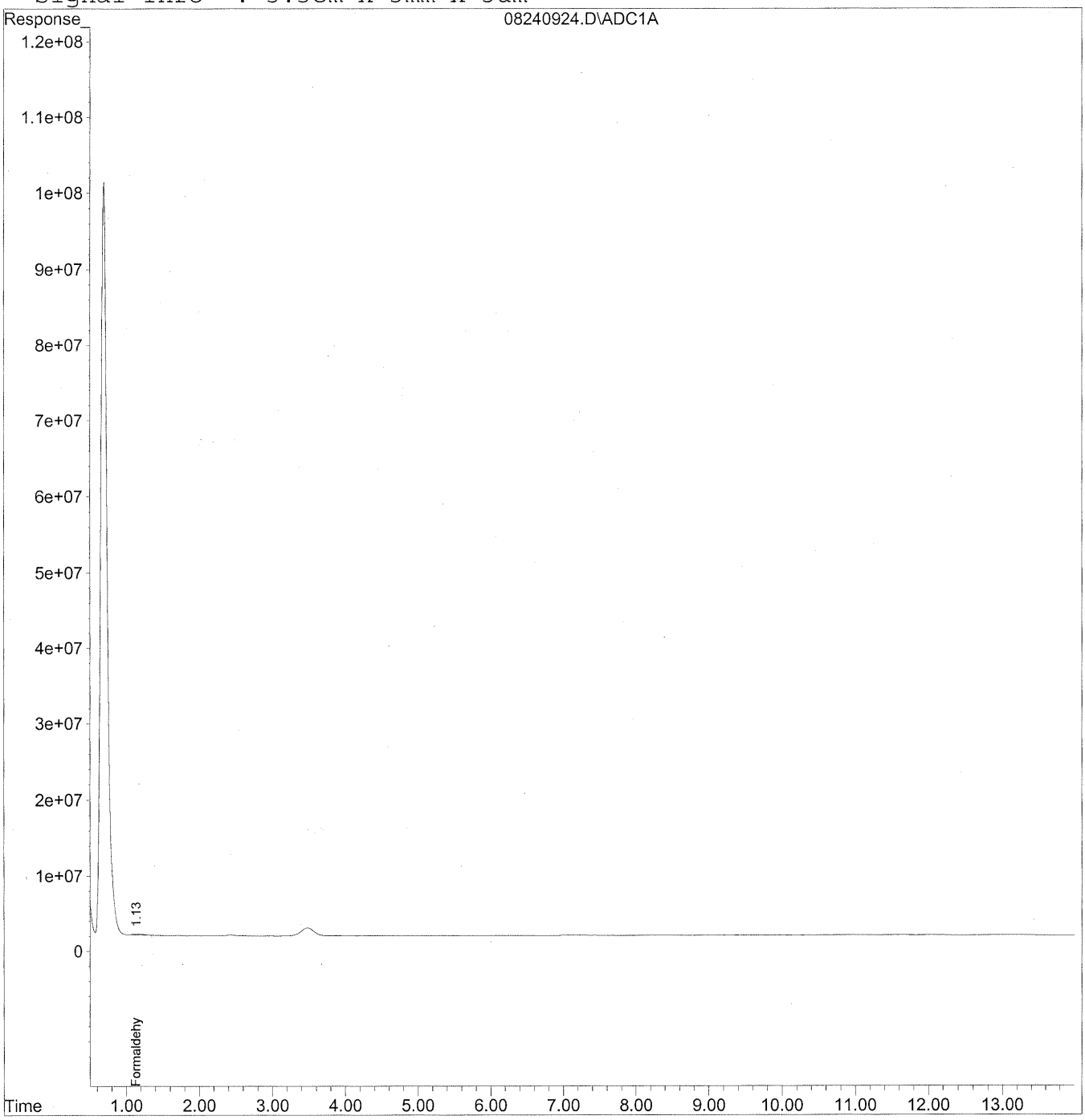
8/31/09

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_08\24\08240924.D Vial: 22
Acq On : 24 Aug 2009 6:16 pm Operator: HC
Sample : P0902912-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 13:32 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\24\08240924.D Vial: 22
Acq On : 24 Aug 2009 6:16 pm Operator: HC
Sample : P0902912-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 13:32 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

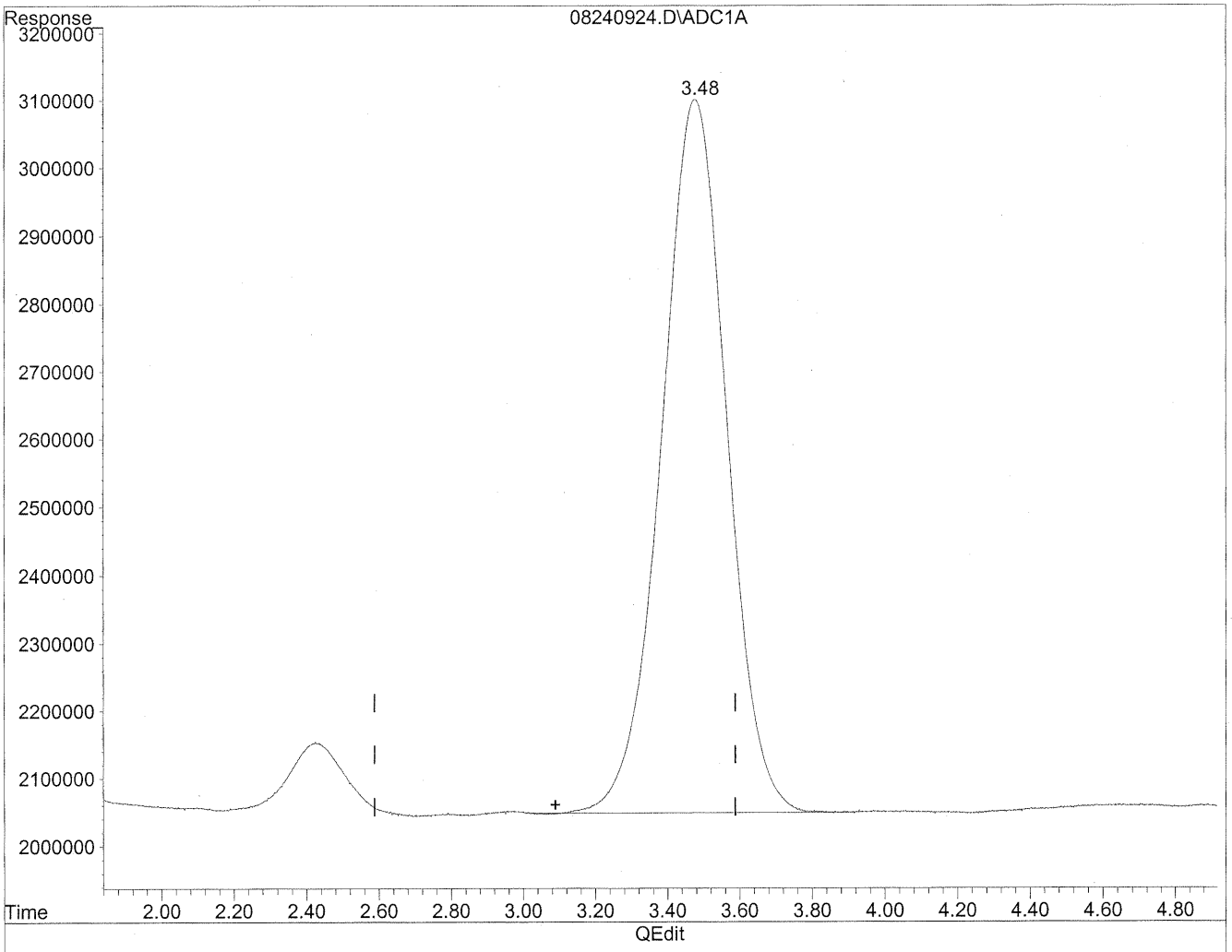
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.13	14014572	76.340	ng/ml
2)	Acetaldehyde	0.00	0	N.D.	ng/ml
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	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

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240924.D Vial: 22
Acq On : 24 Aug 2009 6:16 pm Operator: HC
Sample : P0902912-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 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 12:41:27 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.48min 1238.060ng/ml

response 132095087

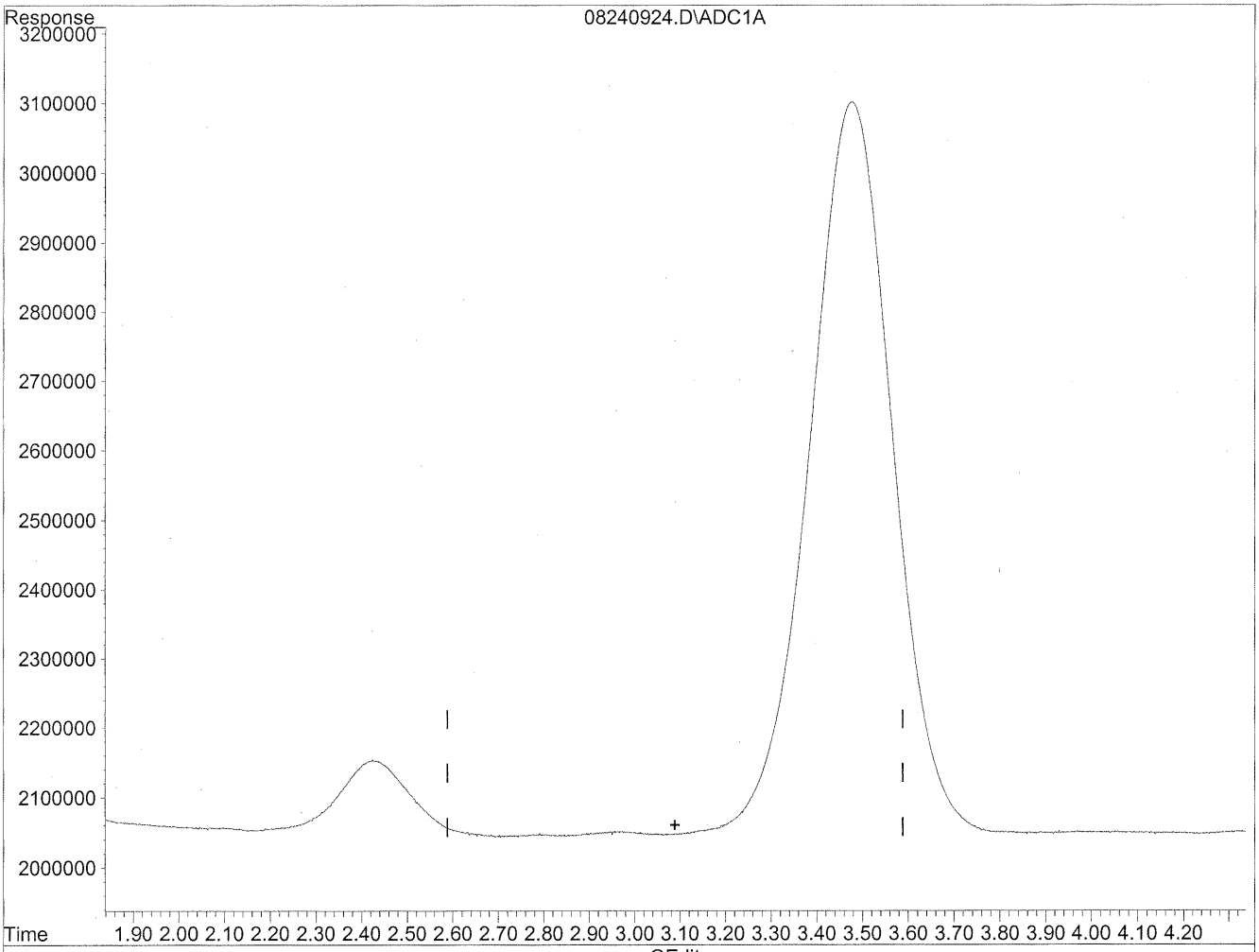
(+) = Expected Retention Time

08240924.D TO110709.M Sat Aug 29 13:32:37 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240924.D Vial: 22
Acq On : 24 Aug 2009 6:16 pm Operator: HC
Sample : P0902912-002 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 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 12:41:27 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC
Shelton
WP

8/31/09

48

(+) = Expected Retention Time
08240924.D TO110709.M Sat Aug 29 13:32:41 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **102512**
Client Project ID: **16512**

CAS Project ID: P0902912
CAS Sample ID: P0902912-003

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: 8/20/09
Date Received: 8/21/09
Date Analyzed: 8/25/09
Desorption Volume: 1.0 ml
Volume Sampled: 108.68 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**5,800**	**53**	0.92	**43**	0.75	
75-07-0	**Acetaldehyde**	**5,500**	**51**	0.92	**28**	0.51	**BT**
123-38-6	**Propionaldehyde**	**530**	**4.9**	0.92	**2.1**	0.39	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.92	ND	0.32	
123-72-8	**Butyraldehyde**	**490**	**4.5**	0.92	**1.5**	0.31	
100-52-7	**Benzaldehyde**	**680**	**6.2**	0.92	**1.4**	0.21	
590-86-3	**Isovaleraldehyde**	**210**	**1.9**	0.92	**0.54**	0.26	
110-62-3	**Valeraldehyde**	**1,400**	**13**	0.92	**3.7**	0.26	
529-20-4	o-Tolualdehyde	< 100	ND	0.92	ND	0.19	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	1.8	ND	0.37	
66-25-1	**n-Hexaldehyde**	**6,300**	**58**	0.92	**14**	0.22	
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	ND	0.92	ND	0.17	

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

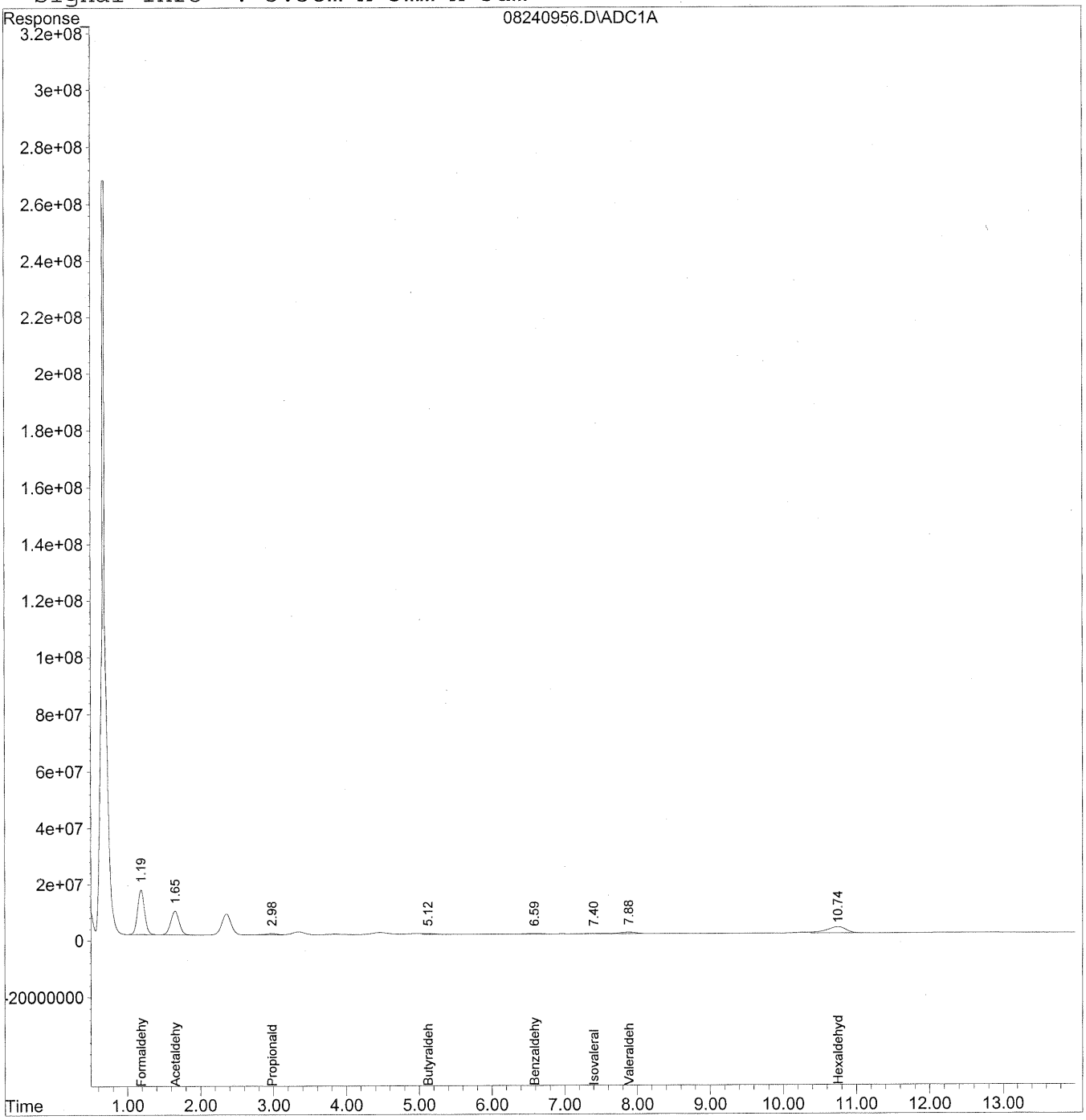
Verified By: Rev Date: 9/2/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 14:53 19109 Quant Results File: TO110709.RES

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



50

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 14:53 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

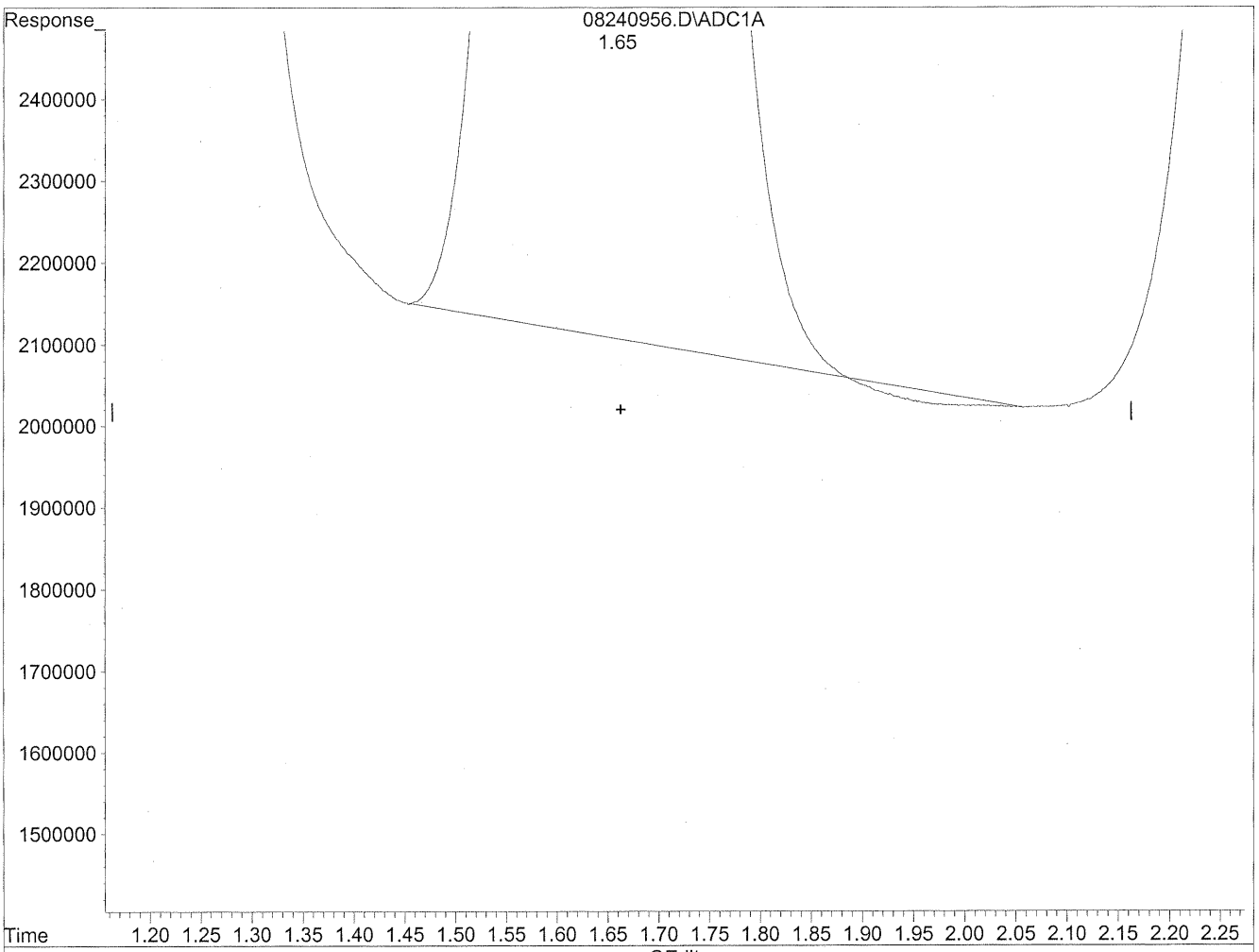
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.19	1064963696	5801.044	ng/ml
2)	Acetaldehyde	1.65	692862713	4941.131	ng/mlm
3)	Propionaldehyde	2.98	56470742	529.272	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	5.12	43272044	489.857	ng/mlm
6)	Benzaldehyde	6.59	44497030	675.535	ng/mlm
7)	Isovaleraldehyde	7.40	16283832	208.097	ng/mlm
8)	Valeraldehyde	7.88	104919625	1427.381	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	10.74	421903917	6264.929	ng/mlm
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/mld

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.65min 4921.427ng/ml

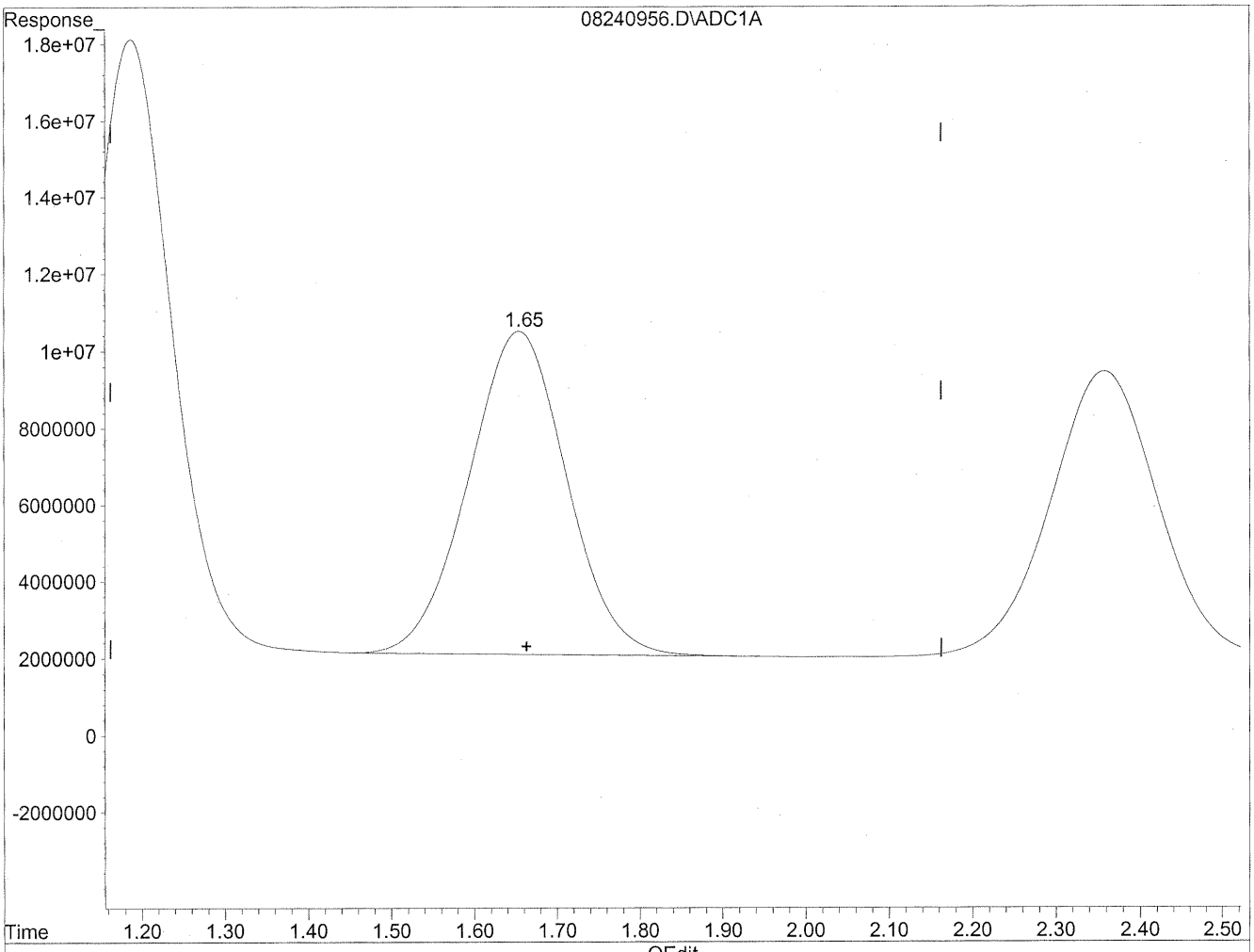
response 690099695

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.65min 4941.131ng/ml m

response 692862713

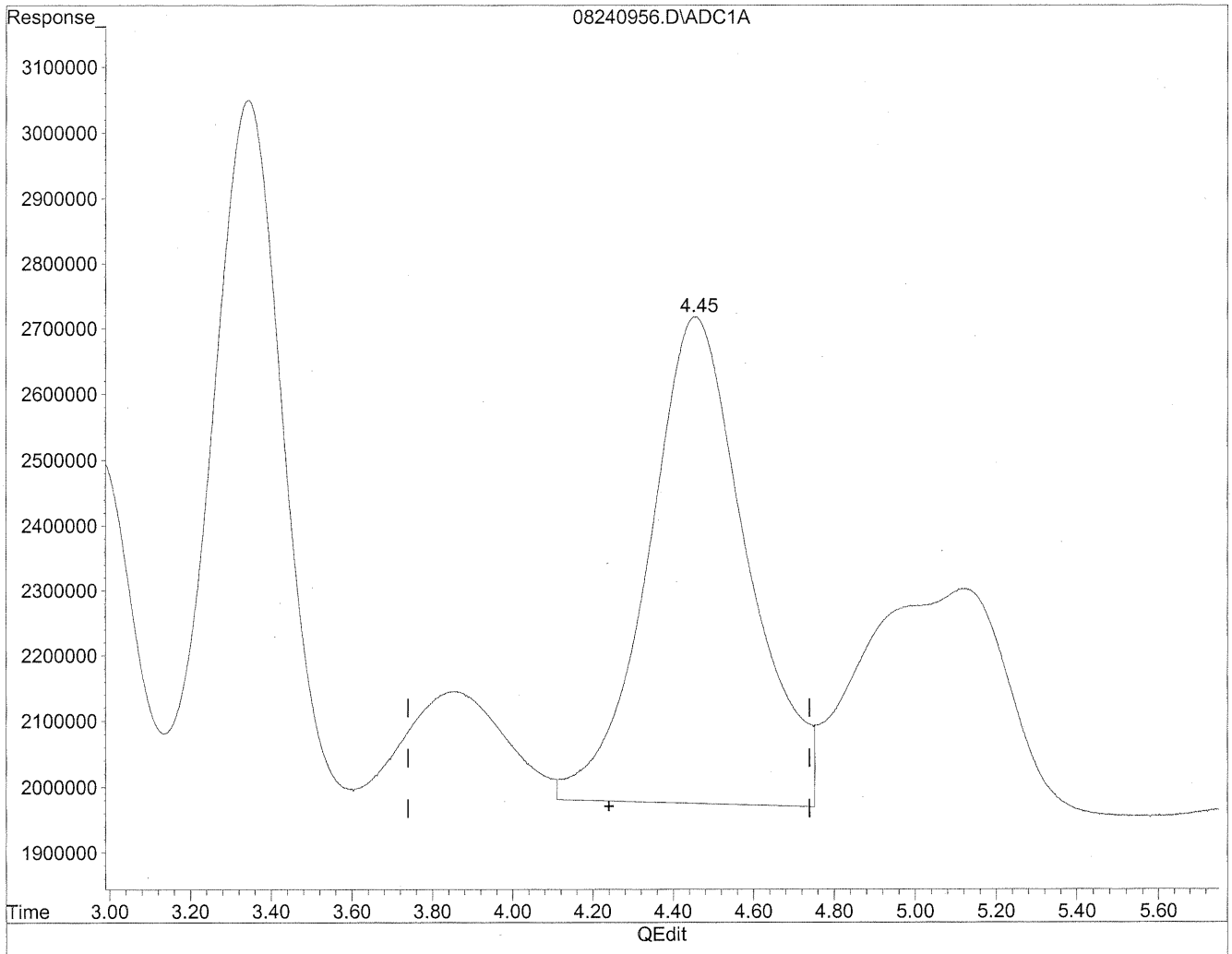
(+) = Expected Retention Time

08240956.D TO110709.M Sat Aug 29 14:49:58 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.46min 1299.715ng/ml

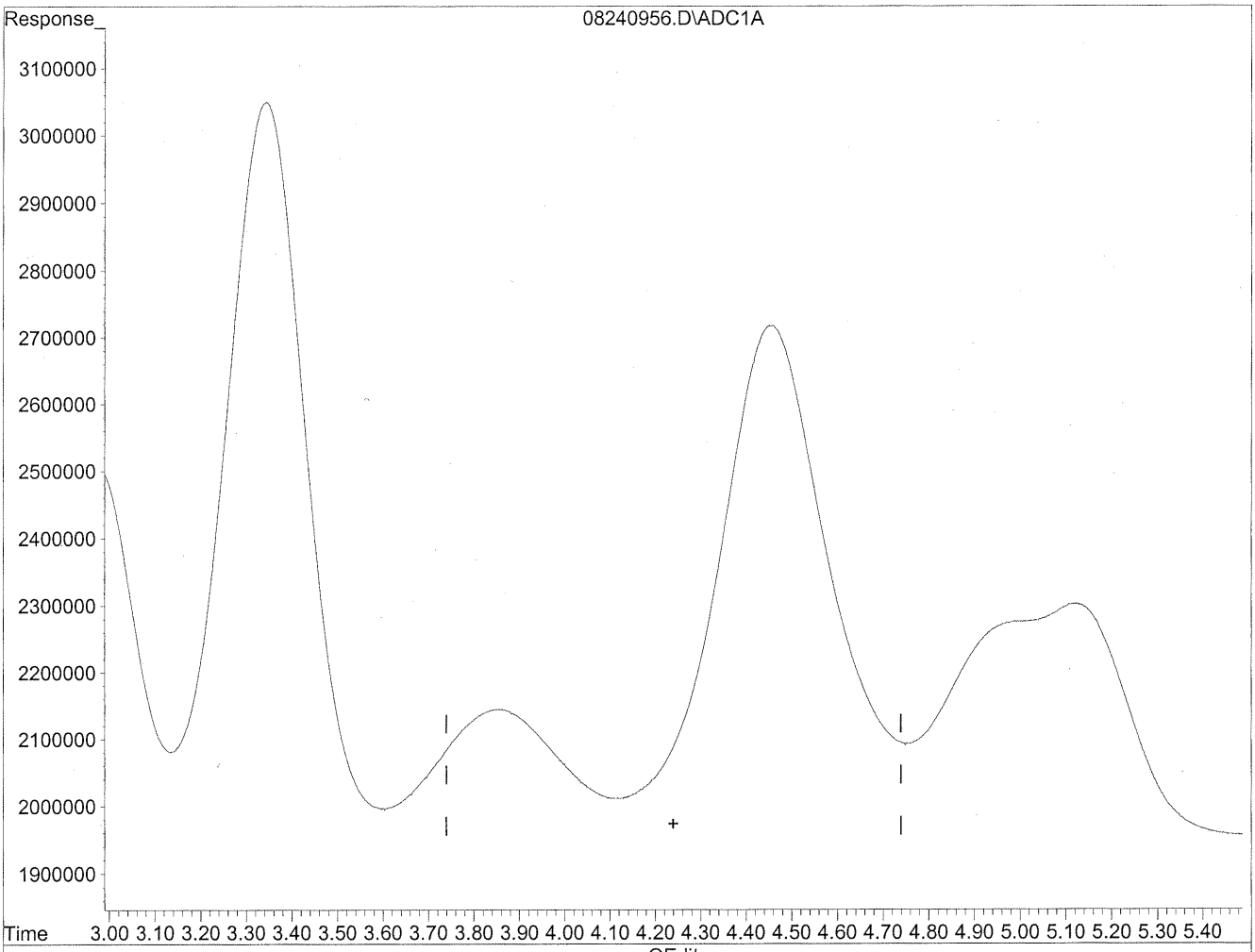
response 126611971

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/31/09 MP

WH 8/31/09

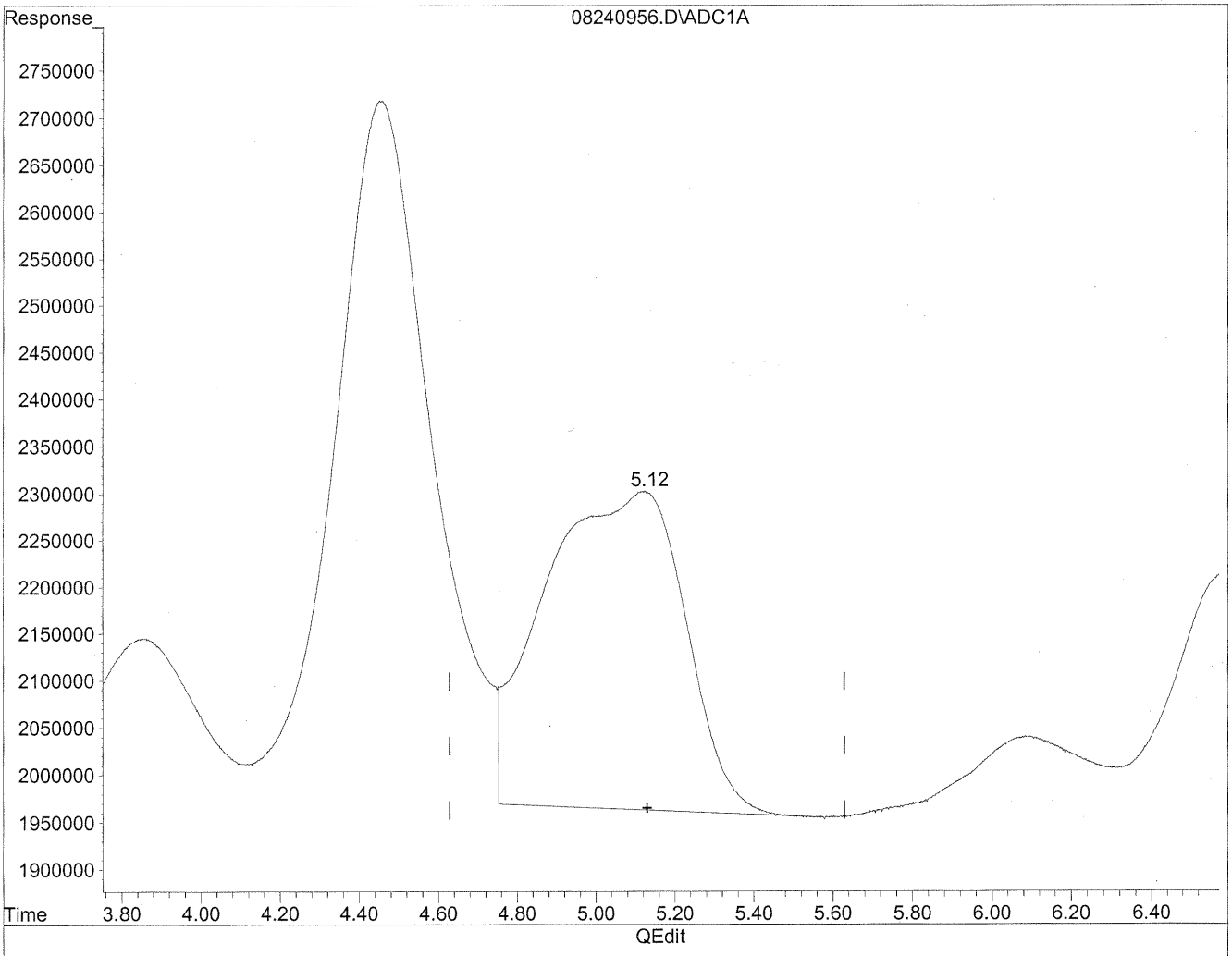
(+) = Expected Retention Time

08240956.D TO110709.M Sat Aug 29 14:50:17 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.12min 949.615ng/ml

response 83885270

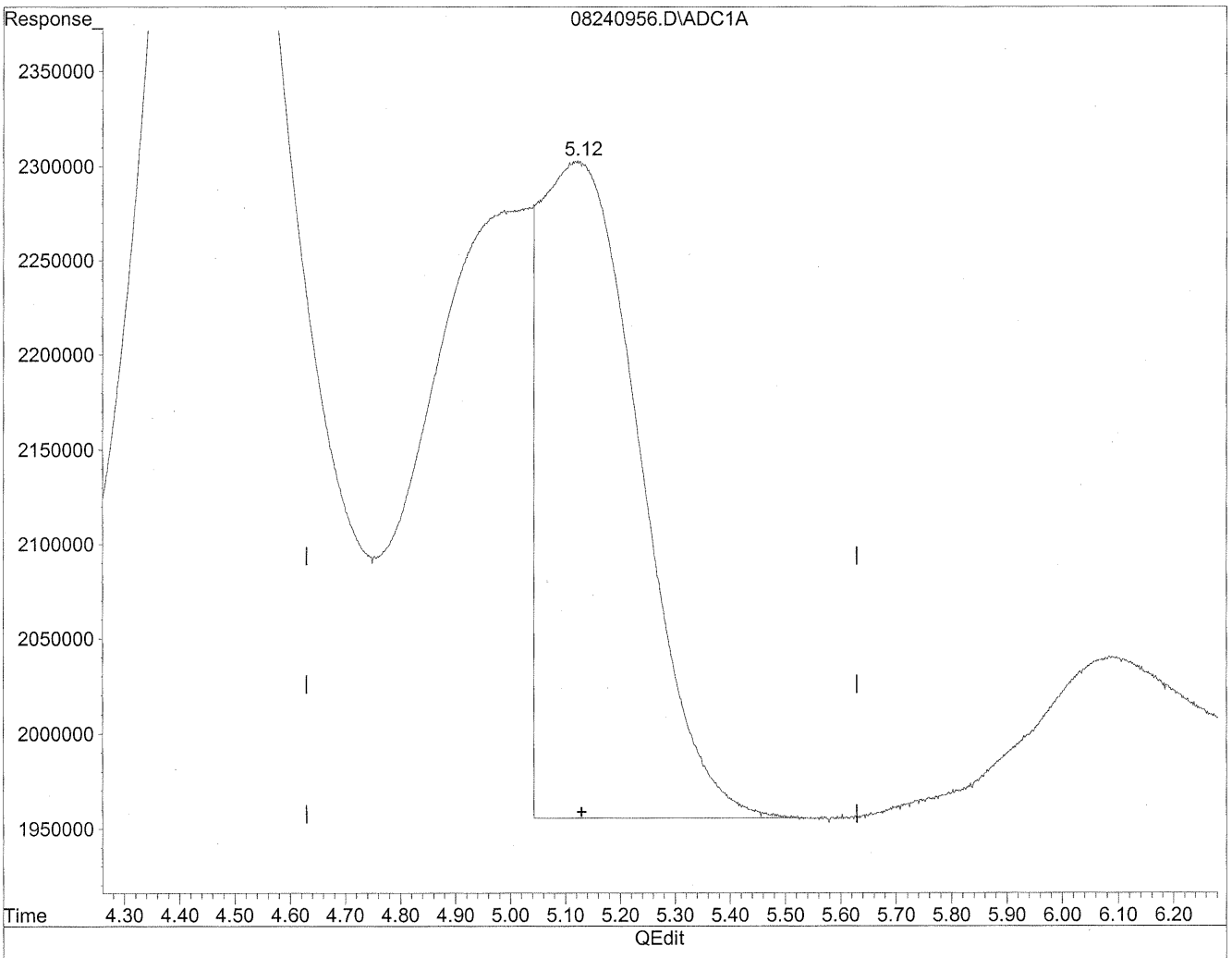
(+) = Expected Retention Time

08240956.D TO110709.M Sat Aug 29 14:50:20 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.12min 489.857ng/ml m

response 43272044

HC 8/29/08 SP

LM 8/31/08

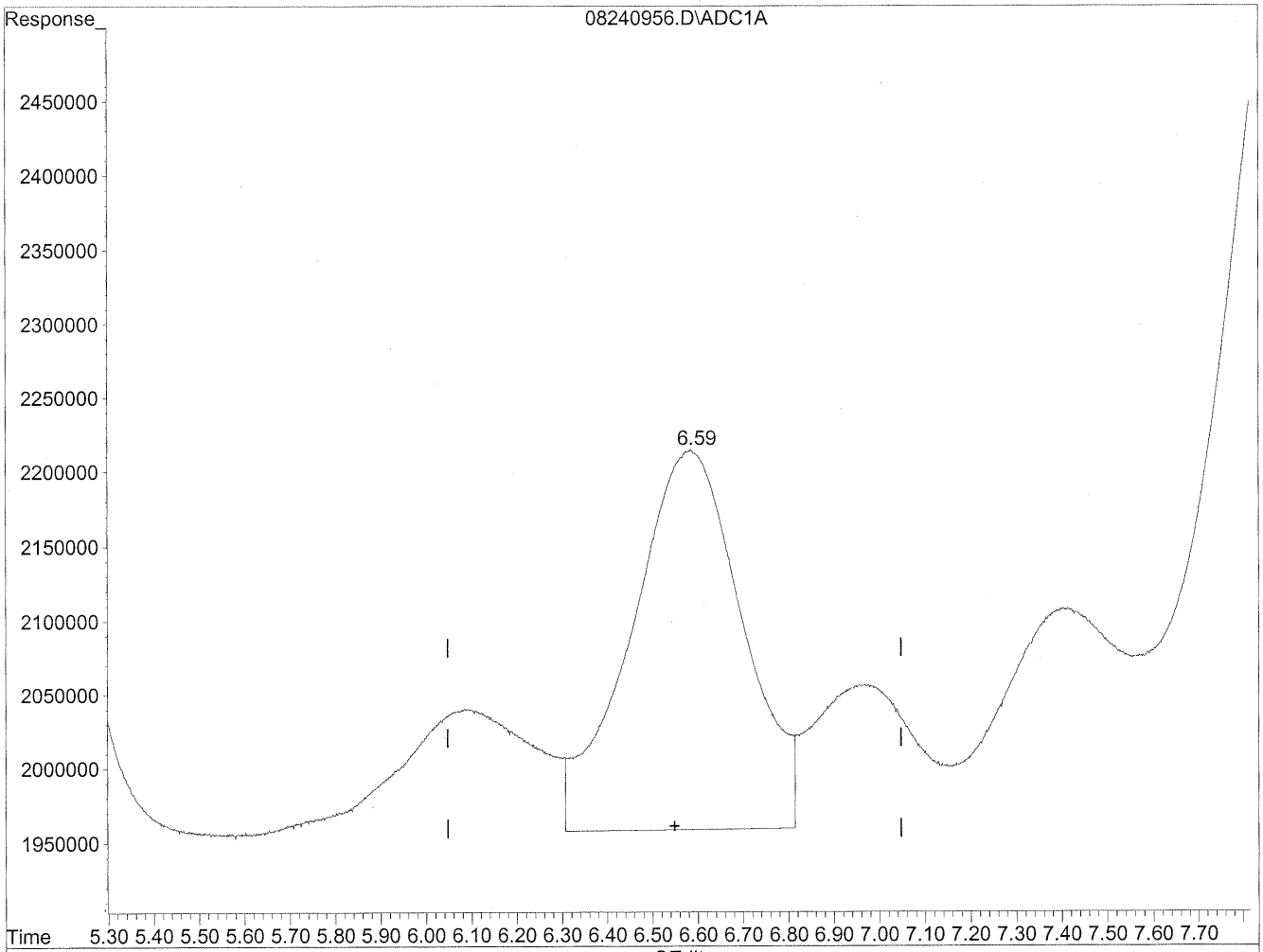
(+) = Expected Retention Time

08240956.D TO110709.M Sat Aug 29 14:50:36 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.58min 663.895ng/ml

response 43730298

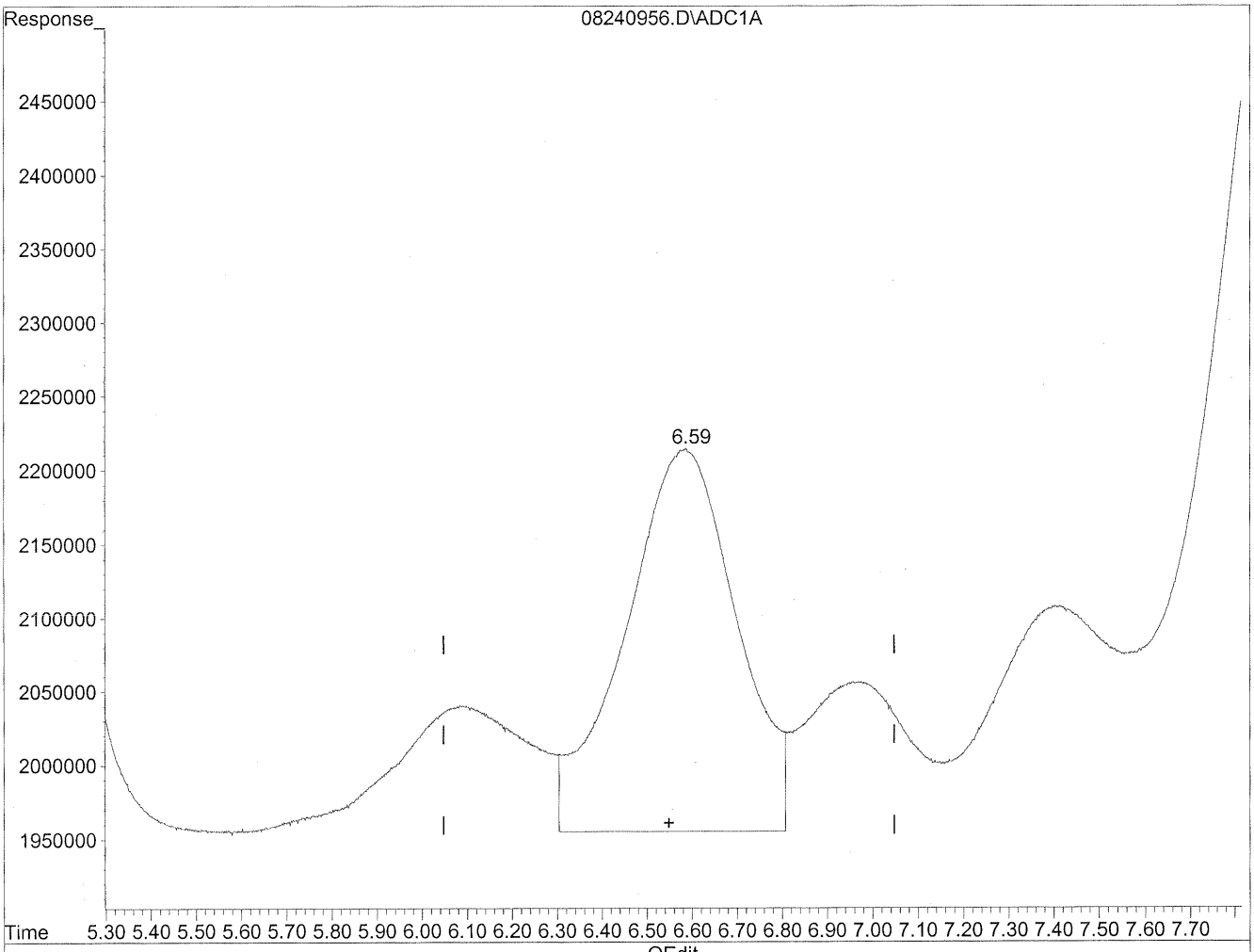
(+) = Expected Retention Time

08240956.D TO110709.M Sat Aug 29 14:50:43 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.59min 675.535ng/ml m

response 44497030

HC 8/29/09

BC

LM 8/31/09

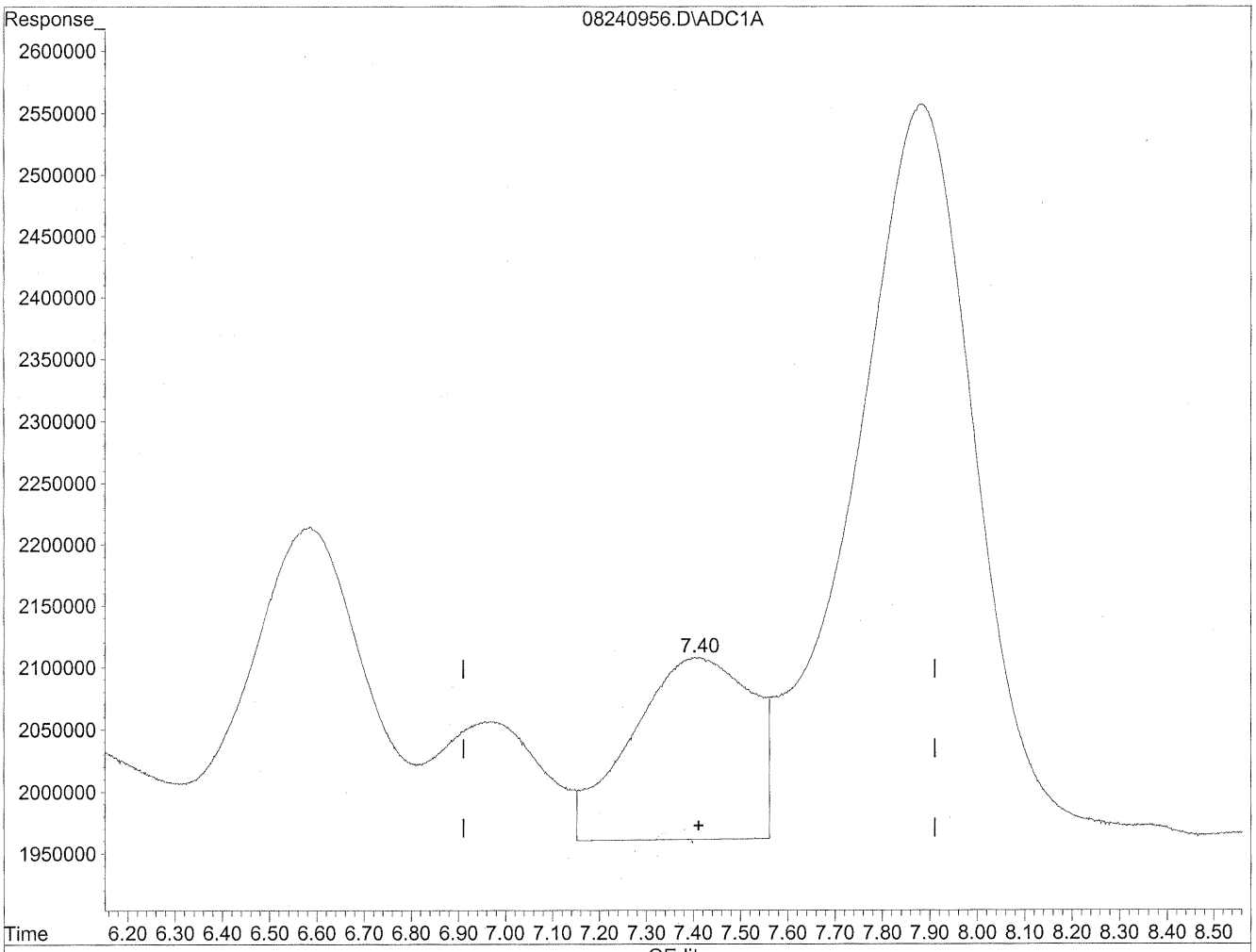
(+) = Expected Retention Time

08240956.D TO110709.M Sat Aug 29 14:51:18 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.41min 333.197ng/ml

response 26073034

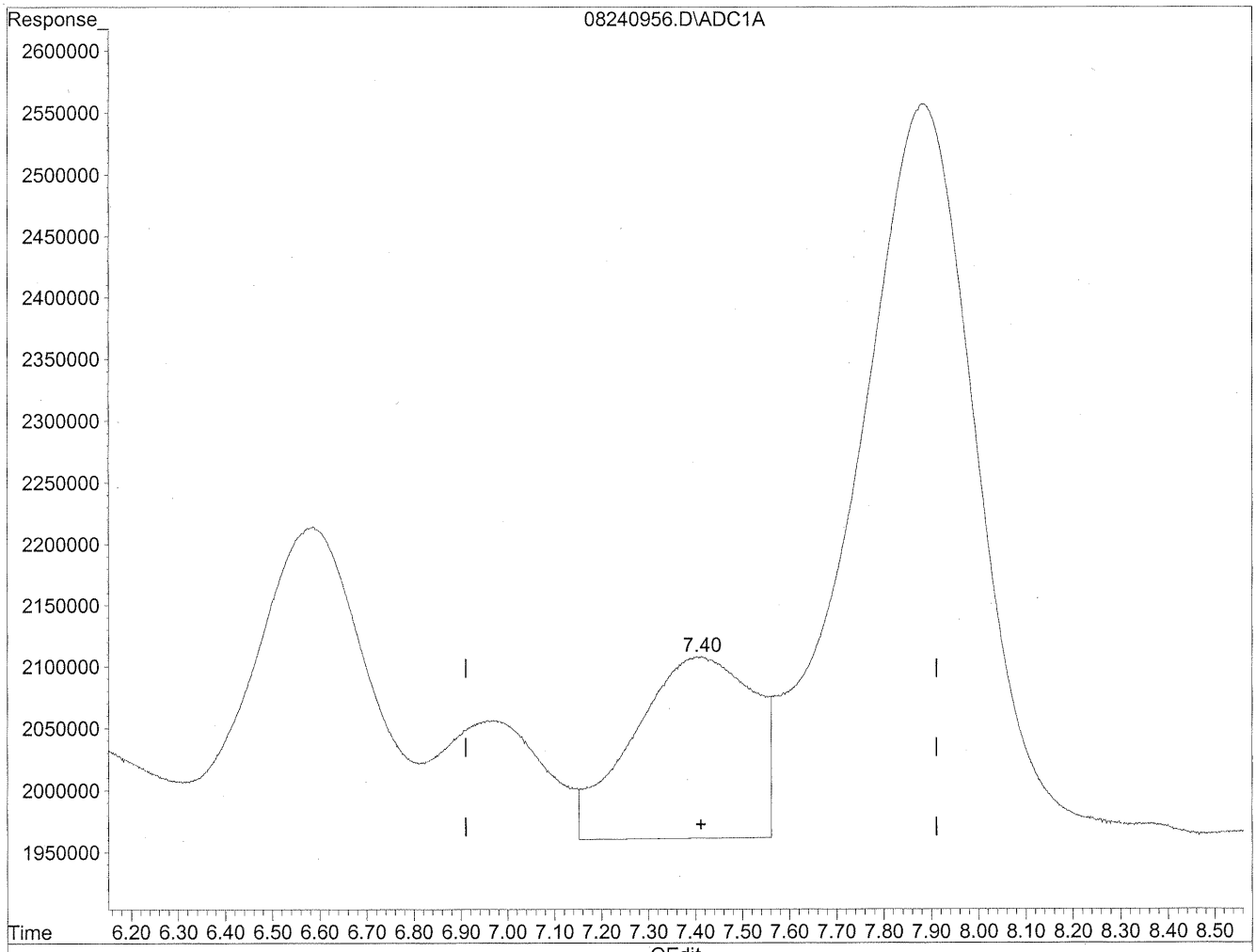
(+) = Expected Retention Time

08240956.D TO110709.M Sat Aug 29 14:51:22 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.41min 333.197ng/ml

response 26073034

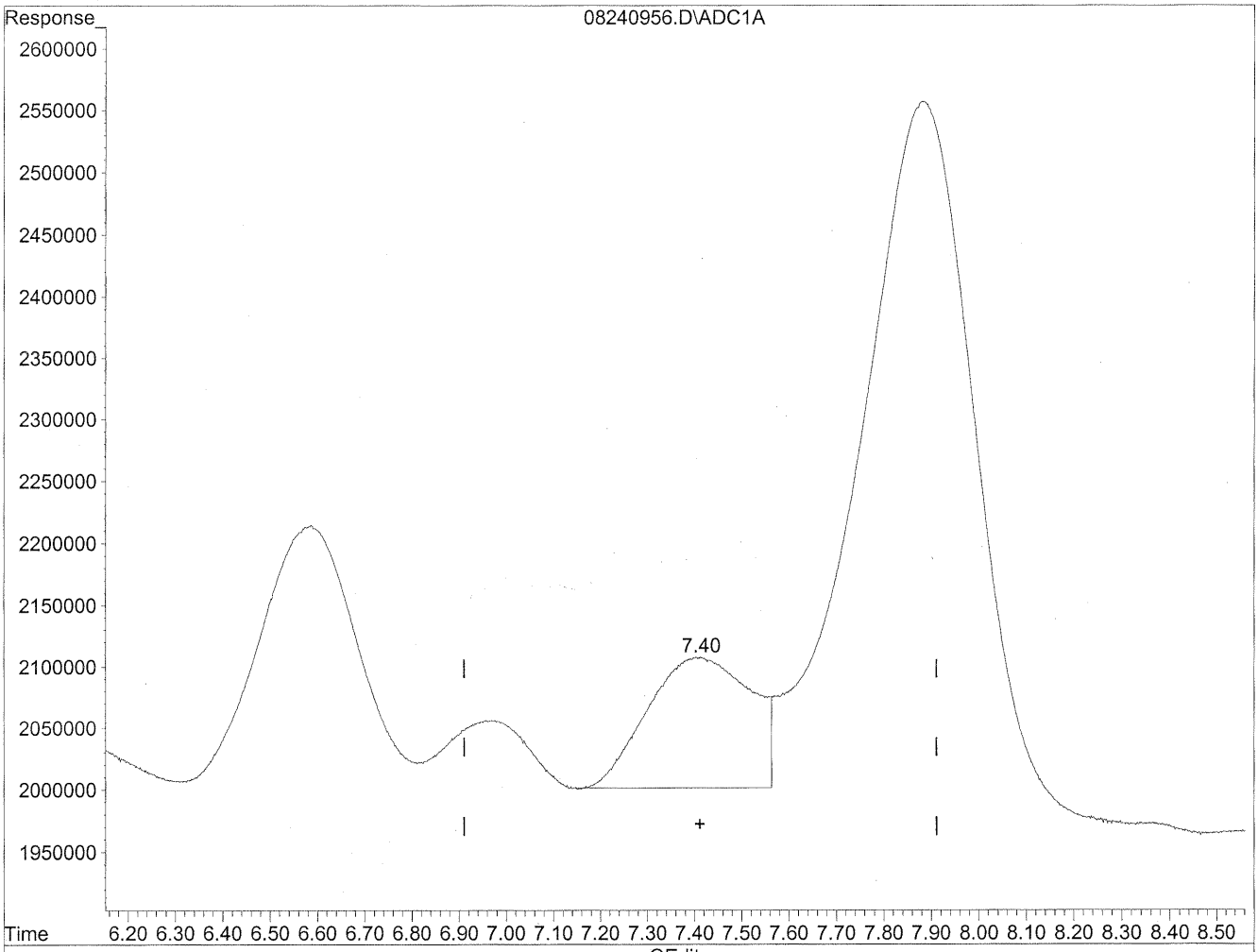
(+) = Expected Retention Time

08240956.D TO110709.M Sat Aug 29 14:51:30 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.40min 208.097ng/ml m

response 16283832

HC station B2

WT 8/31/09

62

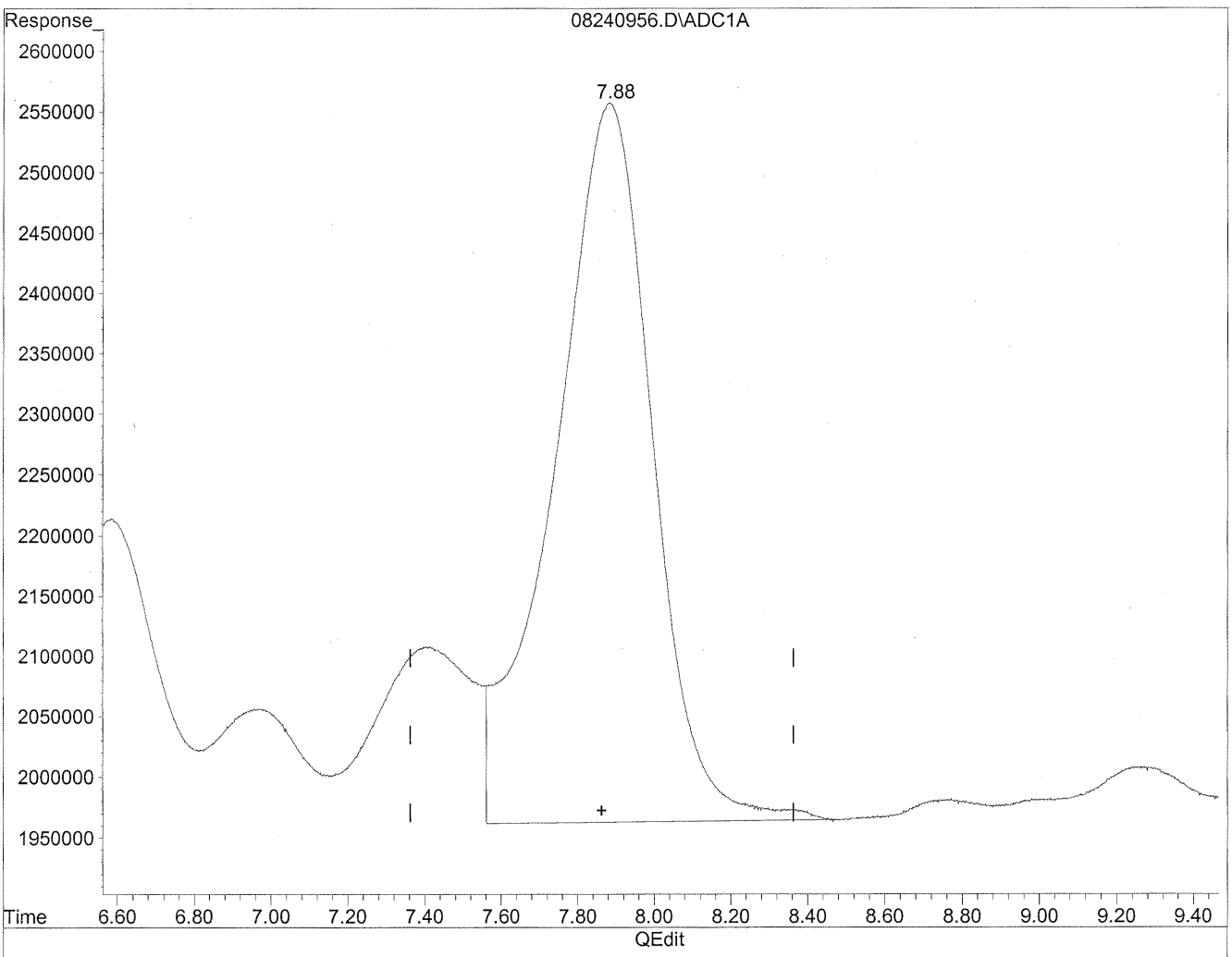
(+) = Expected Retention Time

08240956.D TO110709.M Sat Aug 29 14:51:42 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.88min 1437.269ng/ml

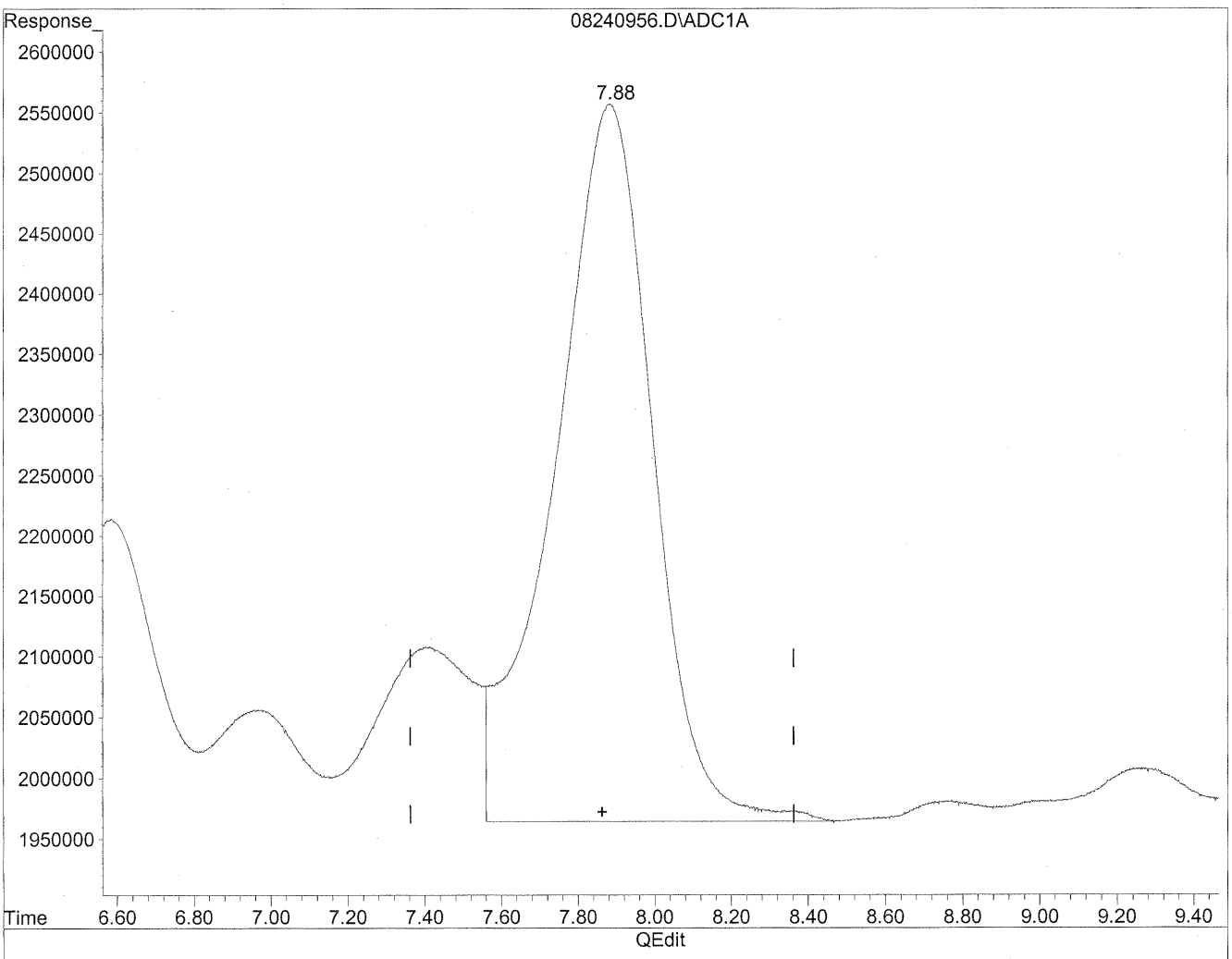
response 105646488

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.88min 1427.381ng/ml m

response 104919625

HC 8/31/09

BC

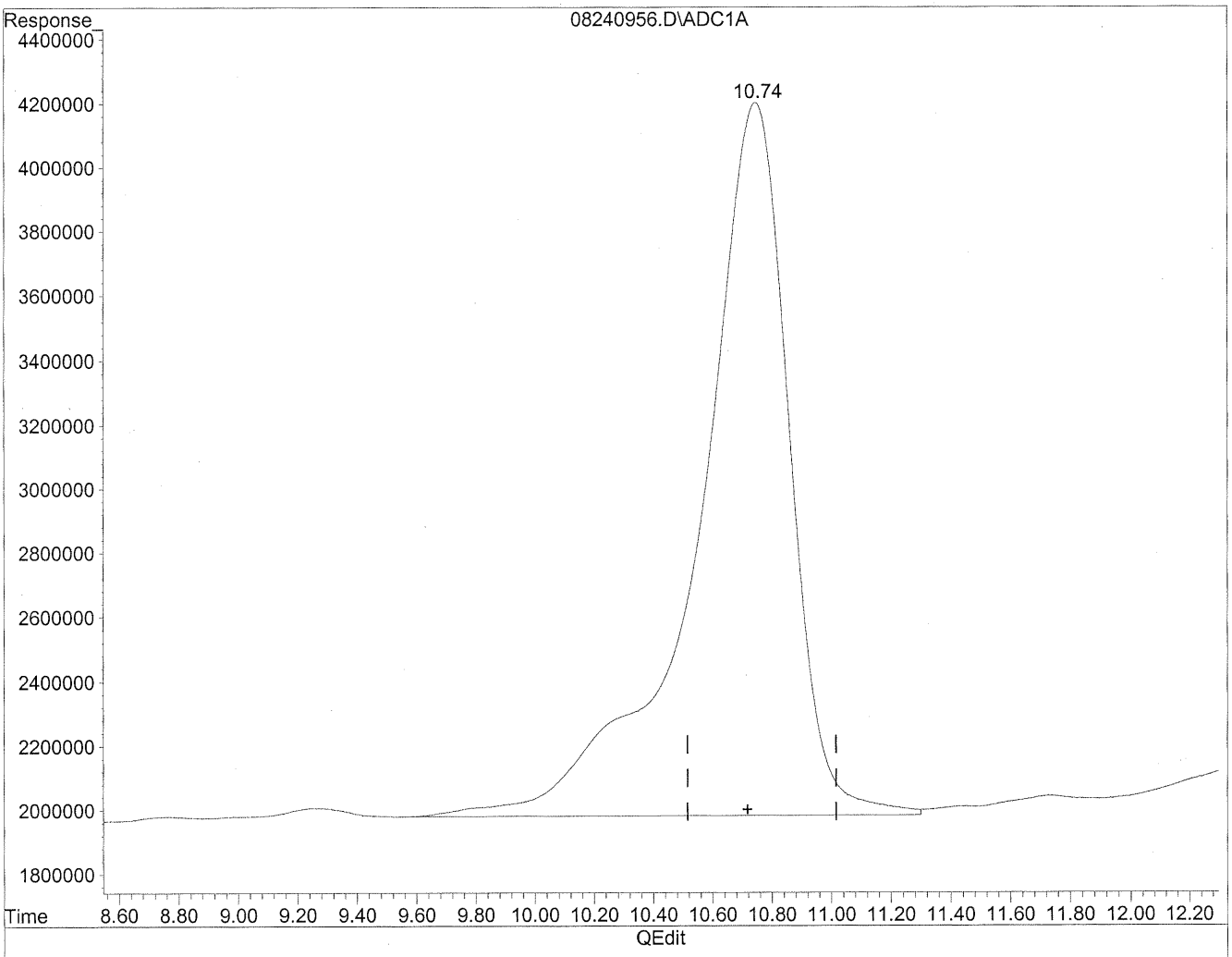
WH 8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.74min 7066.314ng/ml

response 475872163

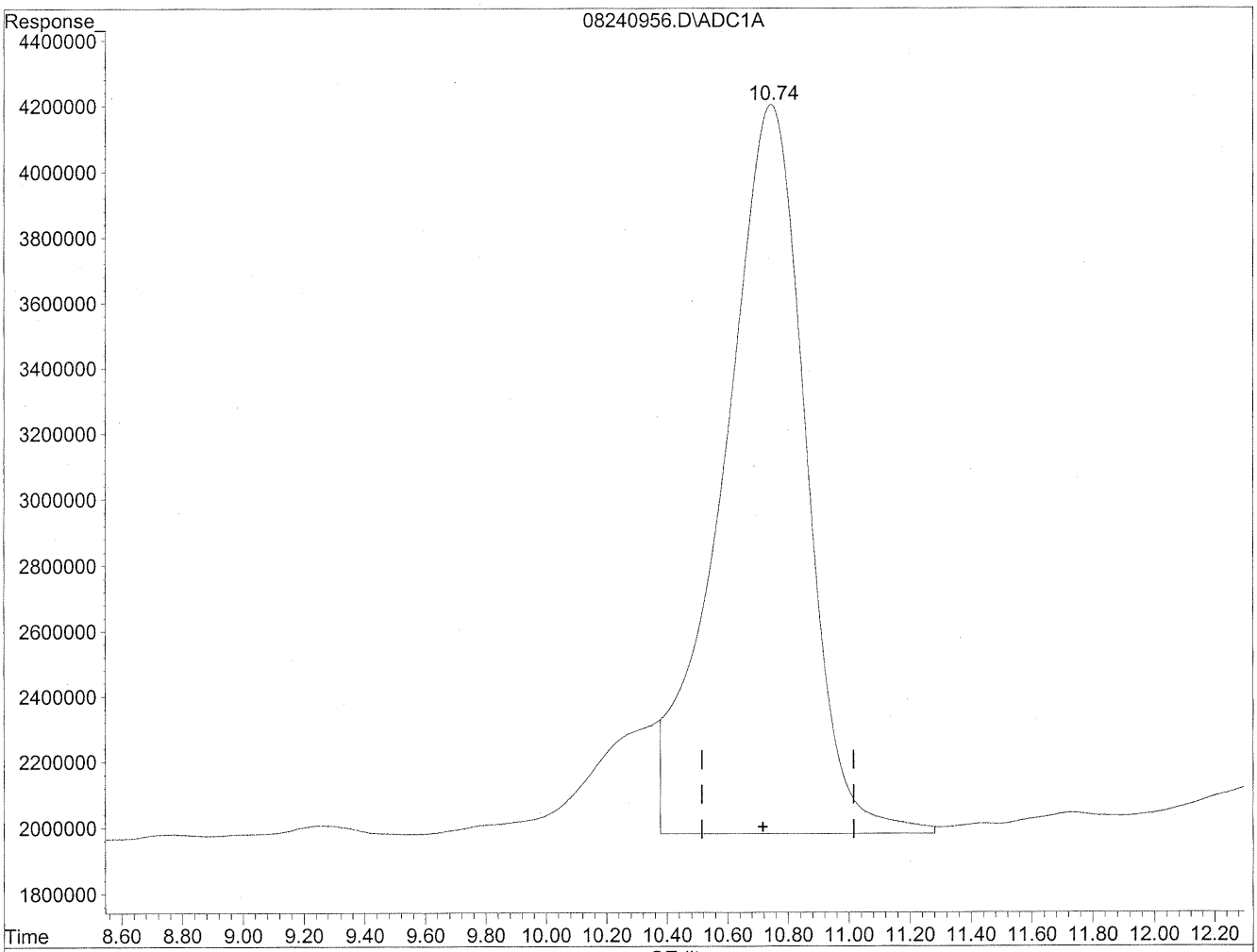
(+) = Expected Retention Time

08240956.D TO110709.M Sat Aug 29 14:52:14 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.74min 6264.929ng/ml m

response 421903917

HC 8/29/09 SH IBC

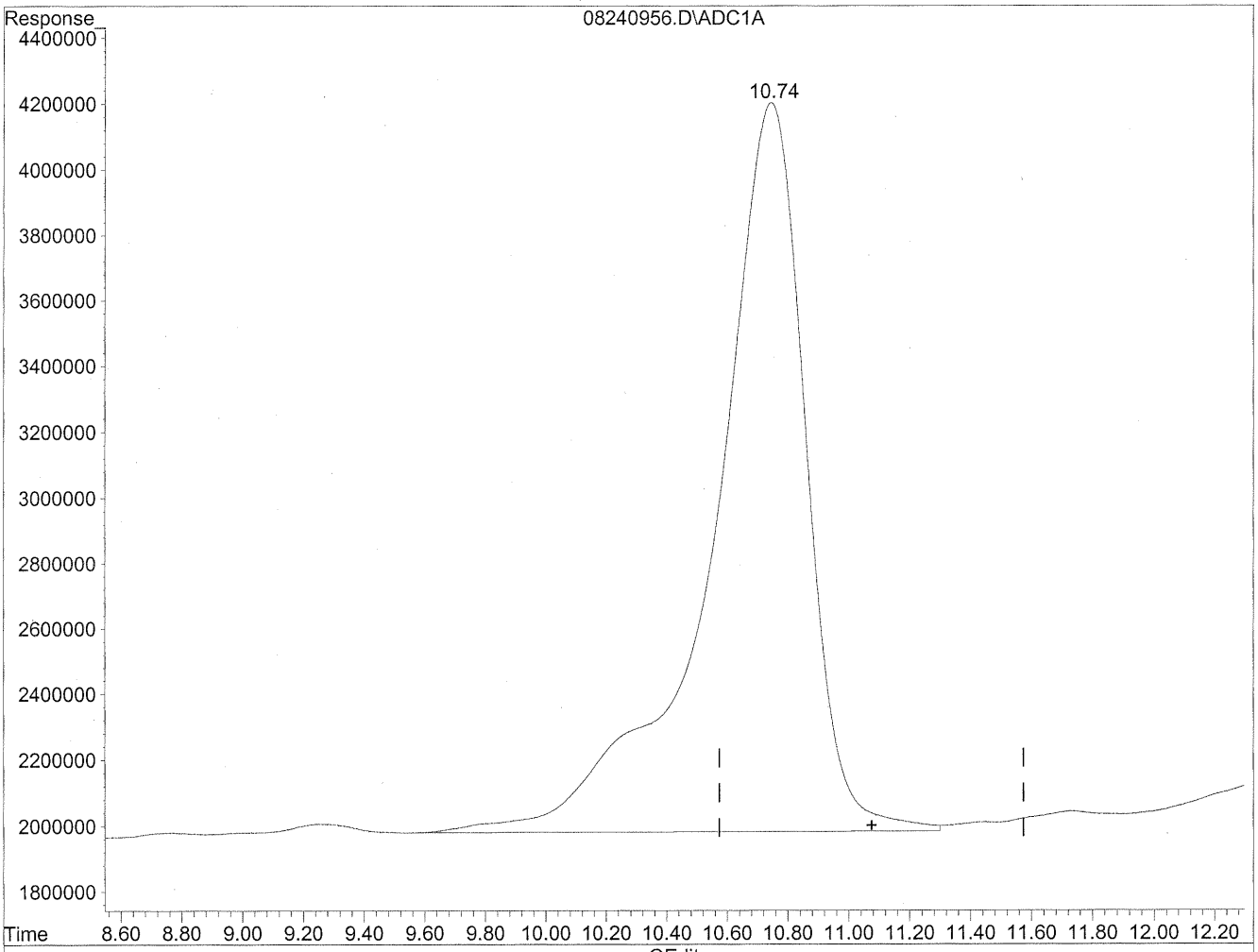
LW 8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.74min 9709.023ng/ml

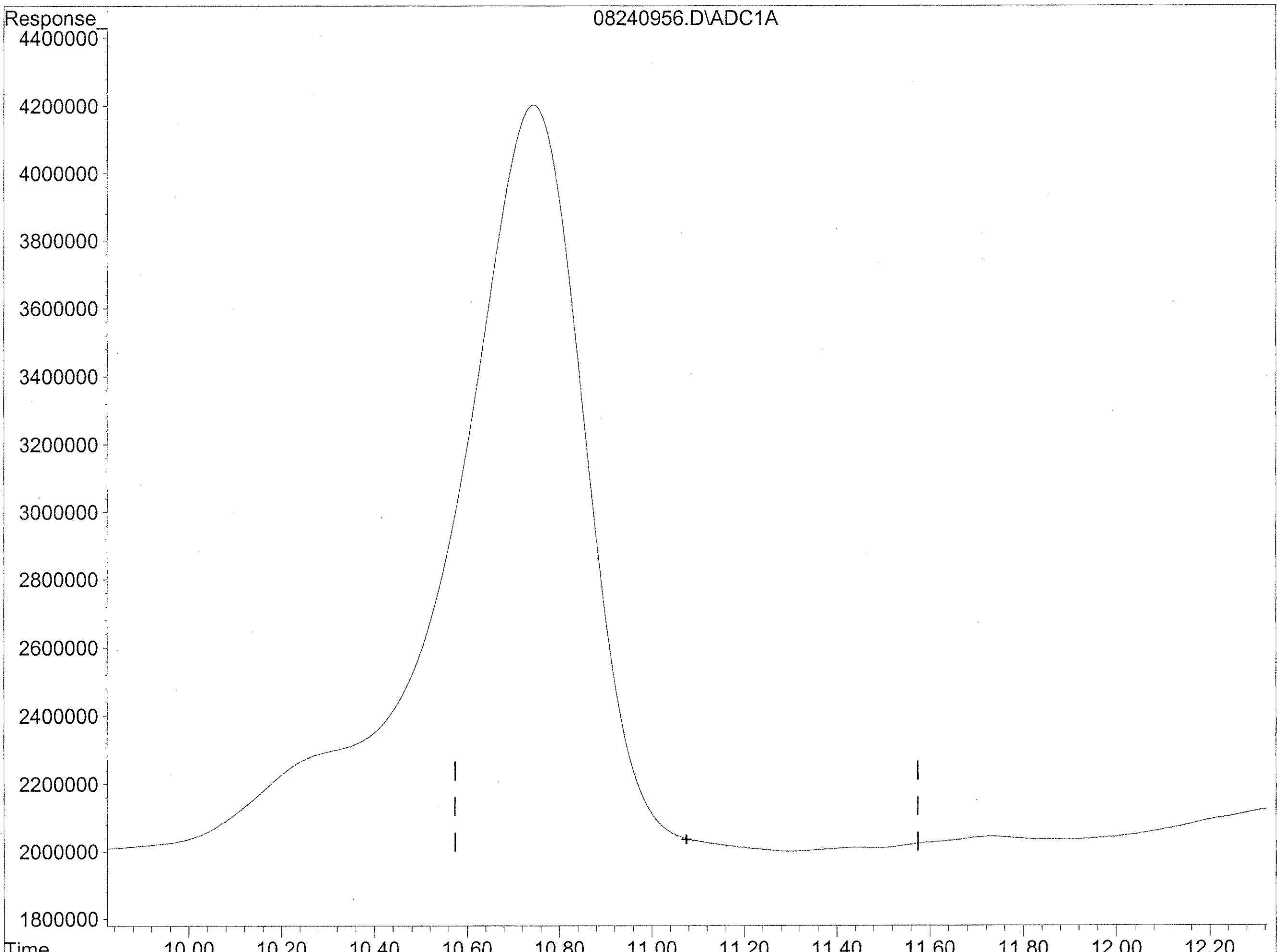
response 475872163

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240956.D Vial: 53
Acq On : 25 Aug 2009 2:17 am Operator: HC
Sample : P0902912-003 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



QEdit

(12) 2,5-Dimethylbenzaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/29/09 MP

WD 8/31/09

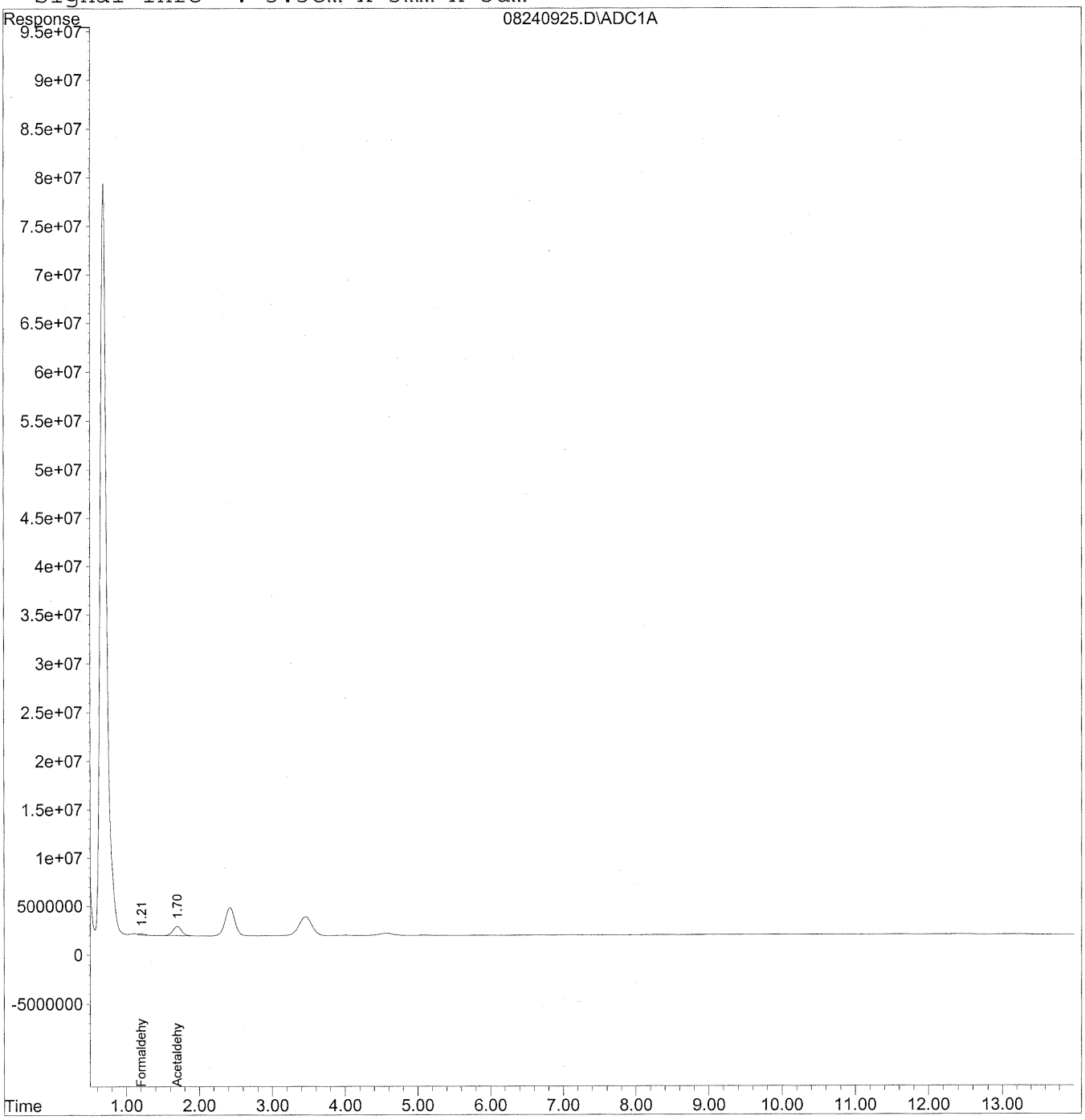
(+) = Expected Retention Time
08240956.D TO110709.M Sat Aug 29 14:53:44 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240925.D Vial: 23
Acq On : 24 Aug 2009 6:31 pm Operator: HC
Sample : P0902912-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 13:37 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 25 10:21:57 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



69

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\24\08240925.D Vial: 23
Acq On : 24 Aug 2009 6:31 pm Operator: HC
Sample : P0902912-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 13:37 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 25 10:21:57 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

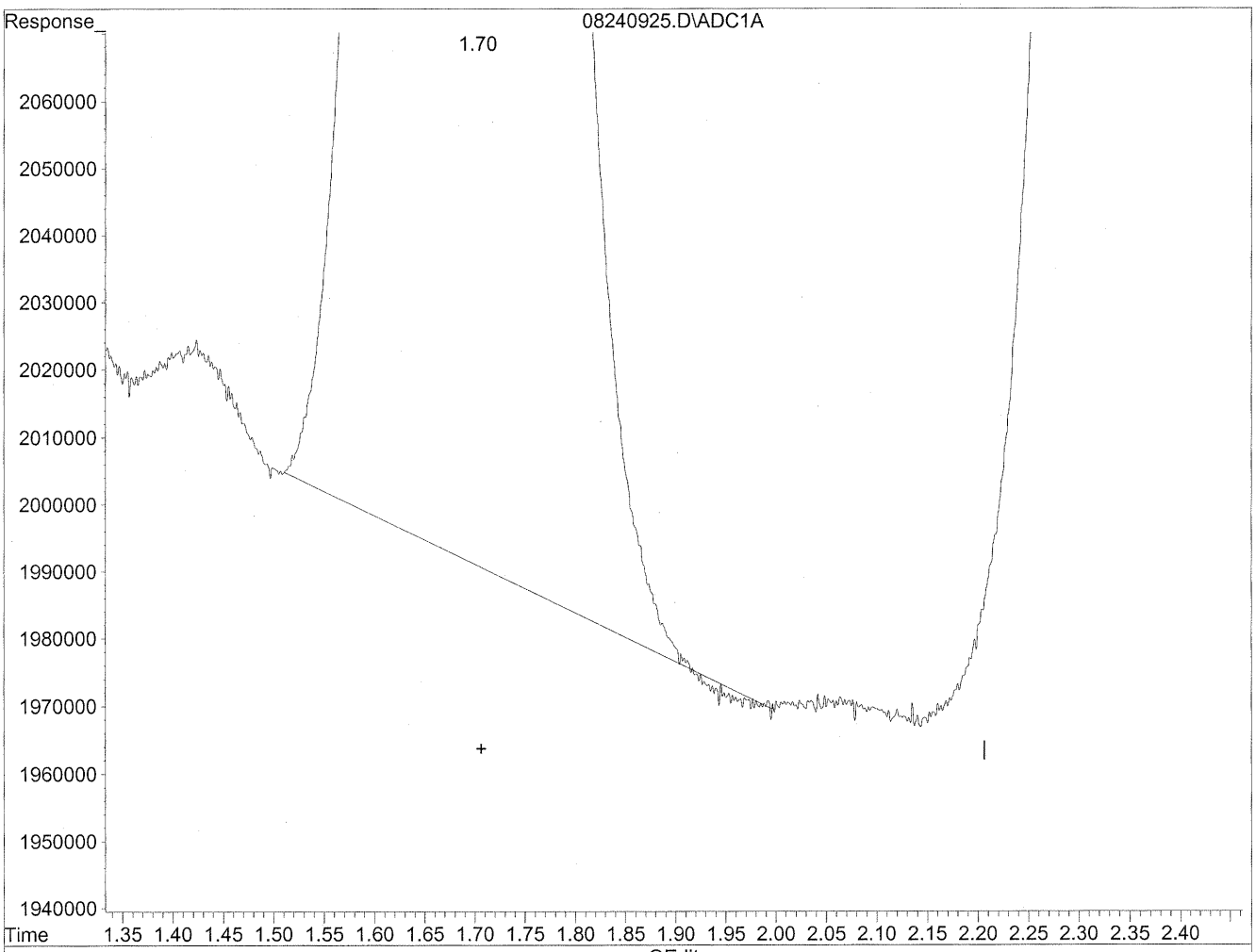
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.20	8305511	45.242	ng/ml
2)	Acetaldehyde	1.70	79639349	567.946	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240925.D Vial: 23
Acq On : 24 Aug 2009 6:31 pm Operator: HC
Sample : P0902912-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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 565.895ng/ml

response 79351805

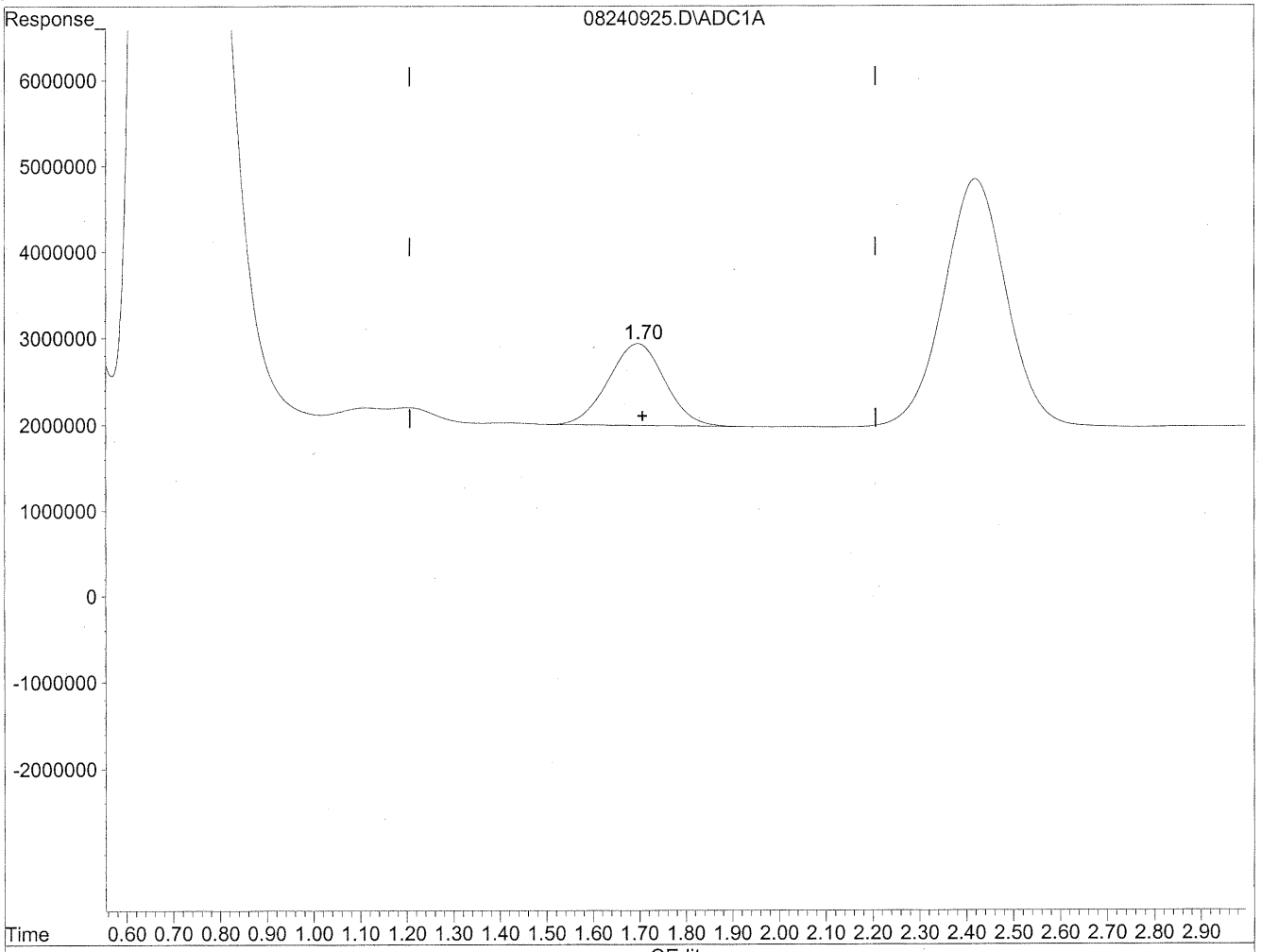
(+) = Expected Retention Time

08240925.D TO110709.M Sat Aug 29 13:36:51 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240925.D Vial: 23
Acq On : 24 Aug 2009 6:31 pm Operator: HC
Sample : P0902912-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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 567.946ng/ml m

response 79639349

tlc 8/29/09

lc

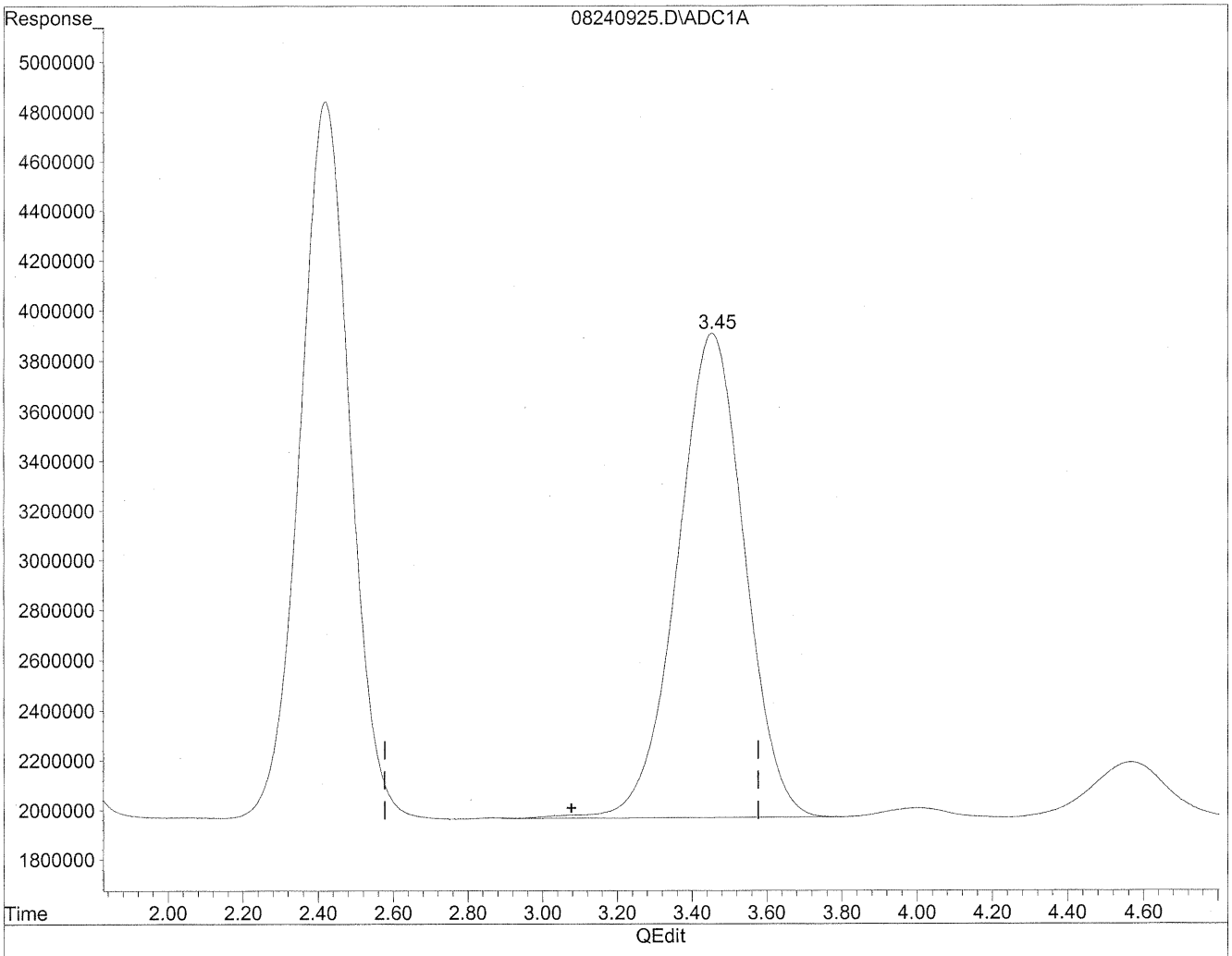
lw 8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240925.D Vial: 23
Acq On : 24 Aug 2009 6:31 pm Operator: HC
Sample : P0902912-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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



(3) Propionaldehyde

3.45min 2278.352ng/ml

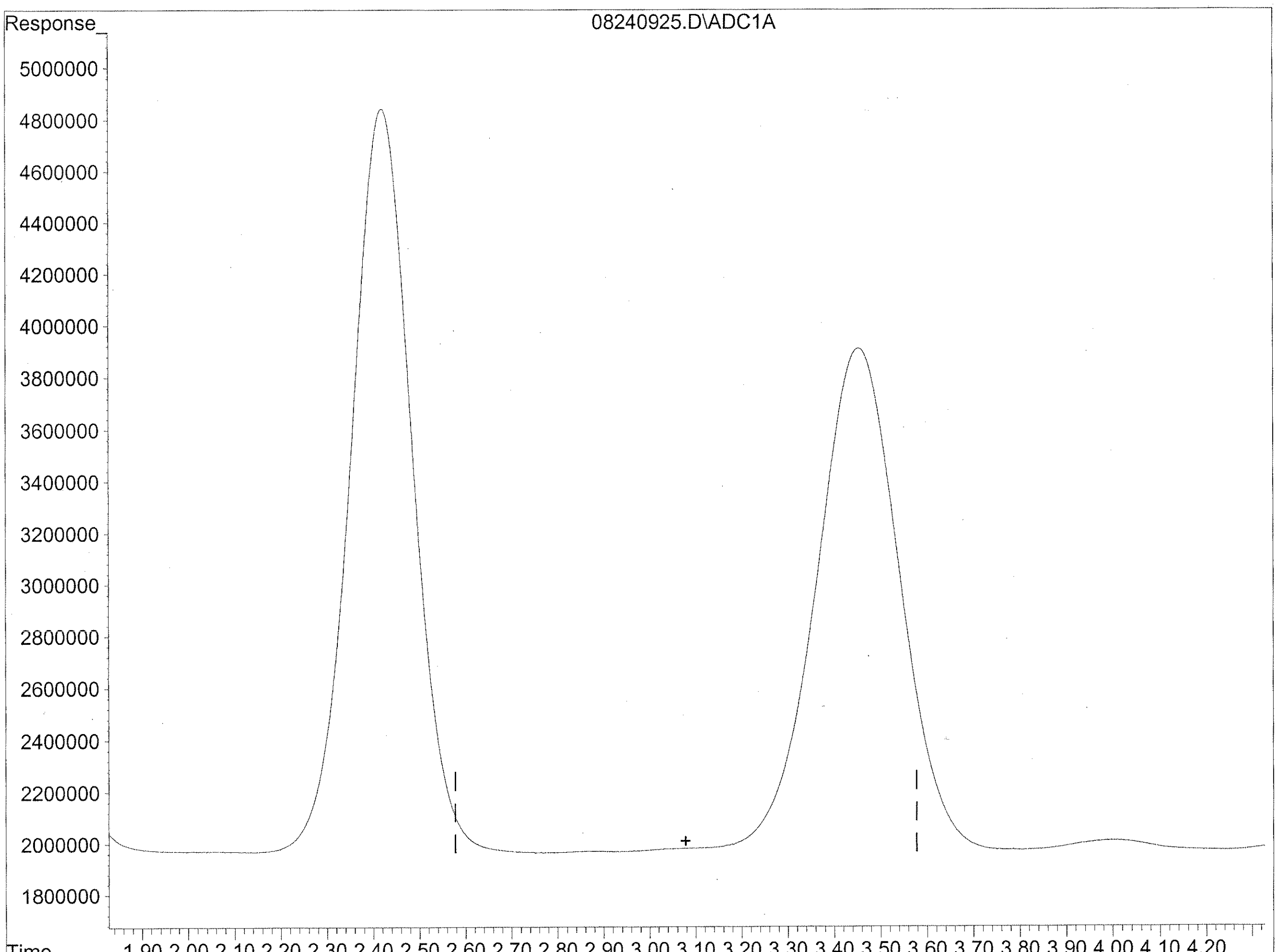
response 243089287

(+) = Expected Retention Time

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2009_08\24\08240925.D       Vial: 23
Acq On    : 24 Aug 2009   6:31 pm                     Operator: HC
Sample    : P0902912-003 back 1.0ml                   Inst    : LC 01
Misc      :                                           Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Aug 25 14:51 19109  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
```



QEdit

(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/29/09 NP

WS 8/31/09

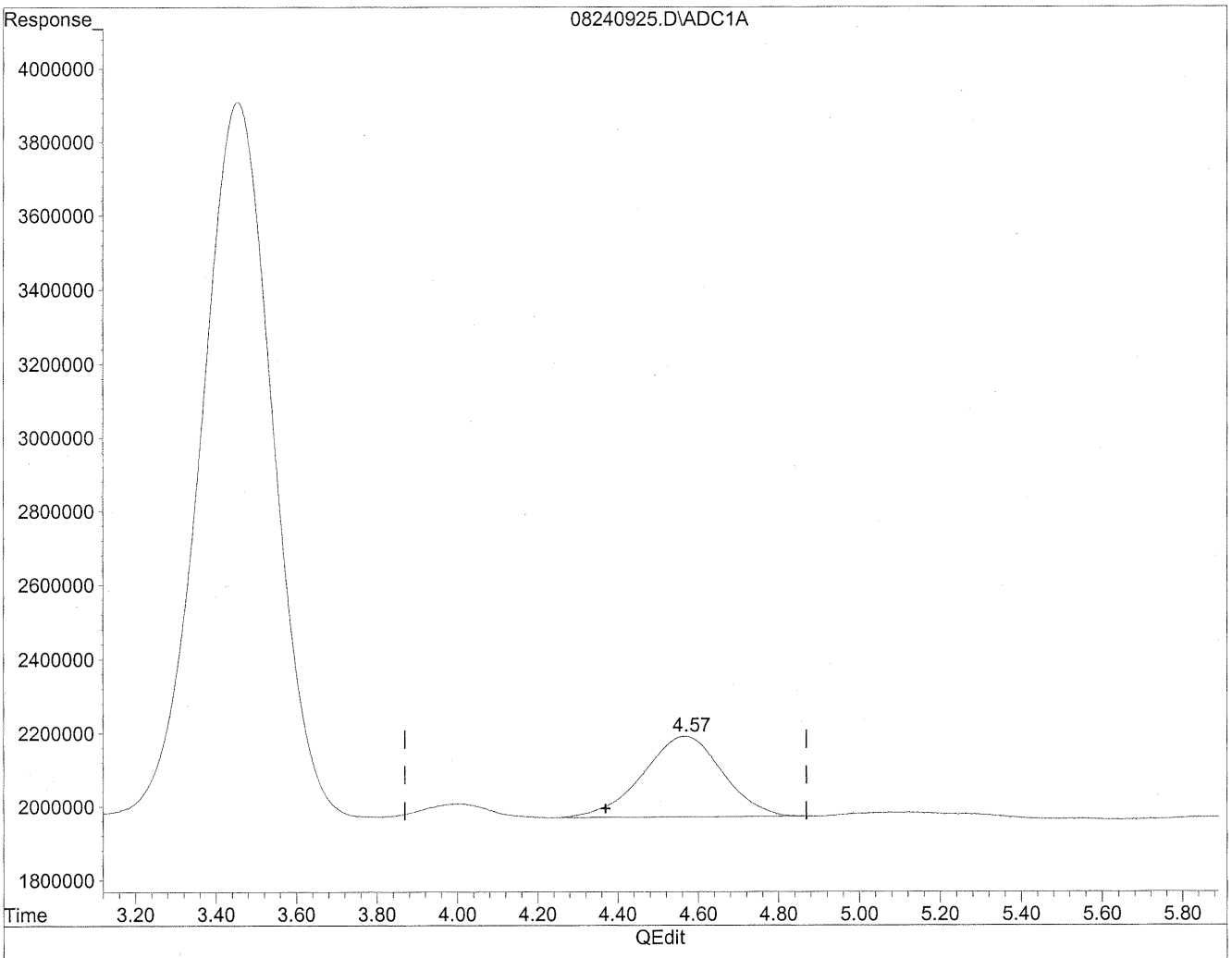
(+) = Expected Retention Time

08240925.D TO110709.M Sat Aug 29 13:37:06 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240925.D Vial: 23
Acq On : 24 Aug 2009 6:31 pm Operator: HC
Sample : P0902912-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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



(4) Crotonaldehyde

4.57min 311.165ng/ml

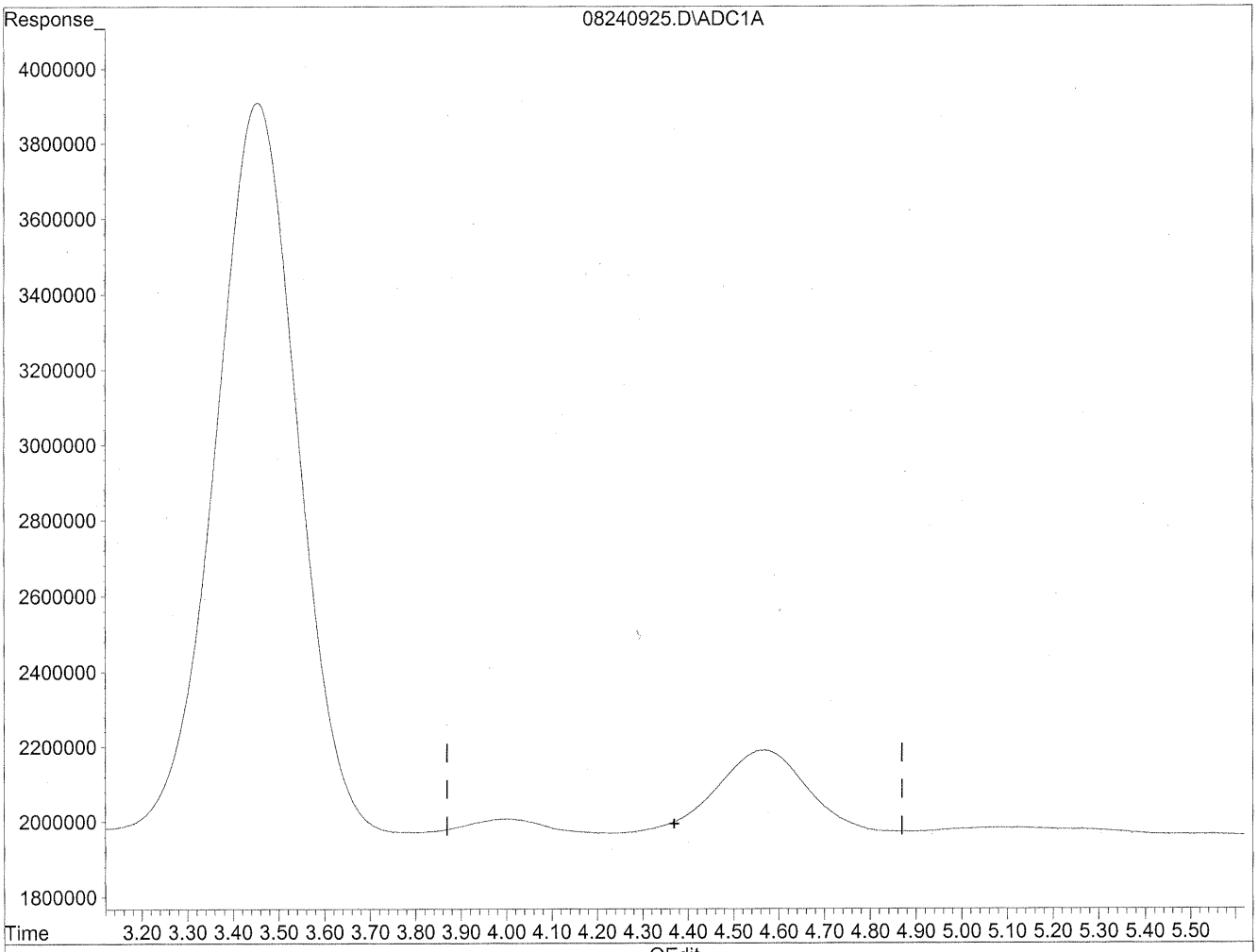
response 30312159

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240925.D Vial: 23
Acq On : 24 Aug 2009 6:31 pm Operator: HC
Sample : P0902912-003 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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



QEdit

(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/29/09 WP

WP 8/31/09

(+) = Expected Retention Time

08240925.D TO110709.M Sat Aug 29 13:37:12 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **102513**
Client Project ID: **16512**

CAS Project ID: P0902912
CAS Sample ID: P0902912-004

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: 8/20/09
Date Received: 8/21/09
Date Analyzed: 8/25/09
Desorption Volume: 1.0 ml
Volume Sampled: 102.96 Liter(s)

CAS #	Compound	Result ng/Sample	Result µg/m³	MRL µg/m³	Result ppbV	MRL ppbV	Data Qualifier
50-00-0	**Formaldehyde**	**5,000**	**49**	0.97	**40**	0.79	
75-07-0	**Acetaldehyde**	**4,900**	**48**	0.97	**27**	0.54	
123-38-6	**Propionaldehyde**	**470**	**4.6**	0.97	**1.9**	0.41	
4170-30-3	Crotonaldehyde, Total	< 100	ND	0.97	ND	0.34	
123-72-8	**Butyraldehyde**	**480**	**4.7**	0.97	**1.6**	0.33	
100-52-7	**Benzaldehyde**	**590**	**5.7**	0.97	**1.3**	0.22	
590-86-3	**Isovaleraldehyde**	**190**	**1.9**	0.97	**0.53**	0.28	
110-62-3	**Valeraldehyde**	**1,300**	**13**	0.97	**3.6**	0.28	
529-20-4	o-Tolualdehyde	< 100	ND	0.97	ND	0.20	
620-23-5							
104-87-0	m,p-Tolualdehyde	< 200	ND	1.9	ND	0.40	
66-25-1	**n-Hexaldehyde**	**5,800**	**56**	0.97	**14**	0.24	
5779-94-2	2,5-Dimethylbenzaldehyde	< 100	ND	0.97	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.

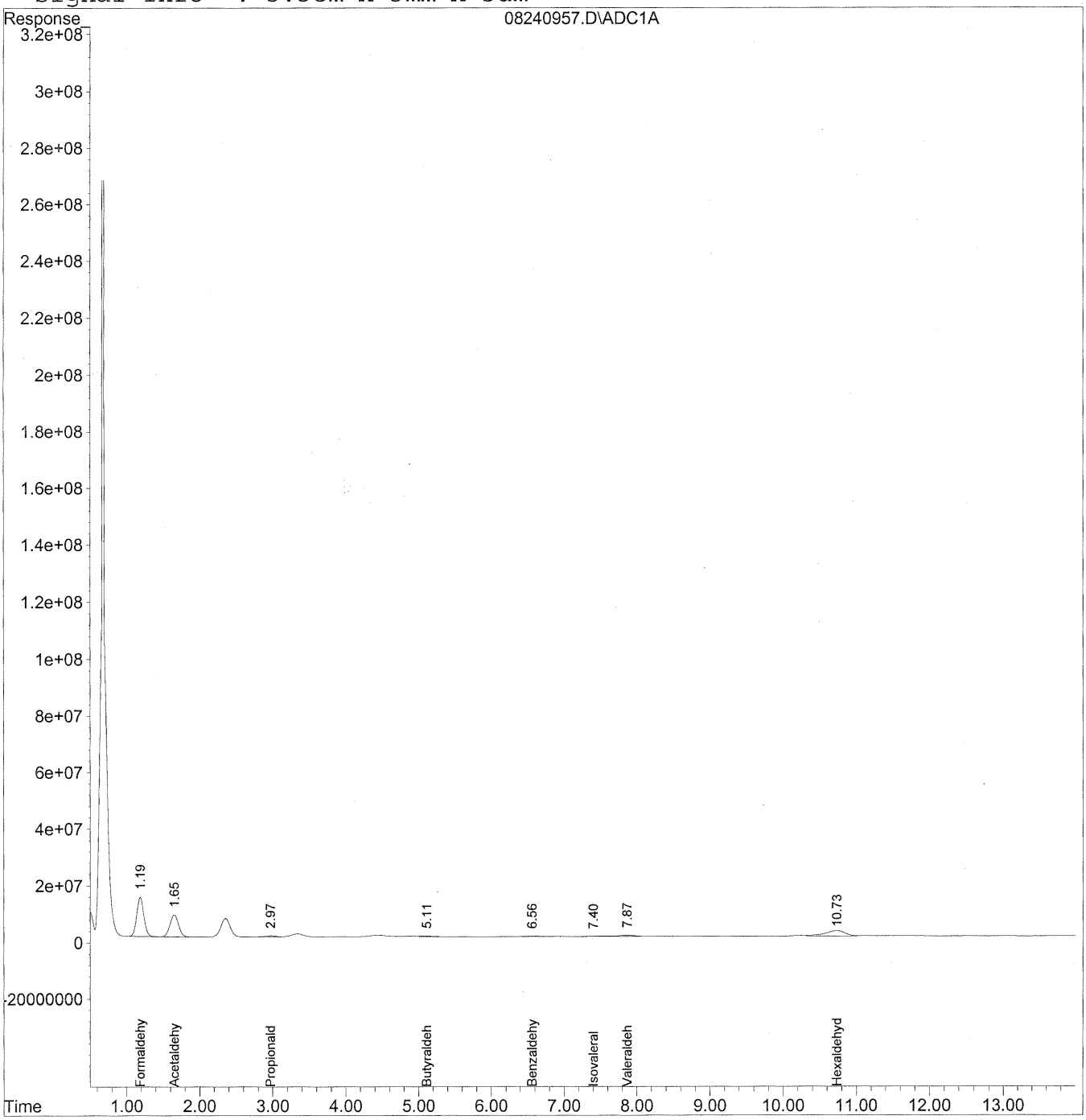
Verified By: Rs Date: 9/2/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 14:58 19109 Quant Results File: TO110709.RES

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



78

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 14:58 19109 Quant Results File: TO110709.RES

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

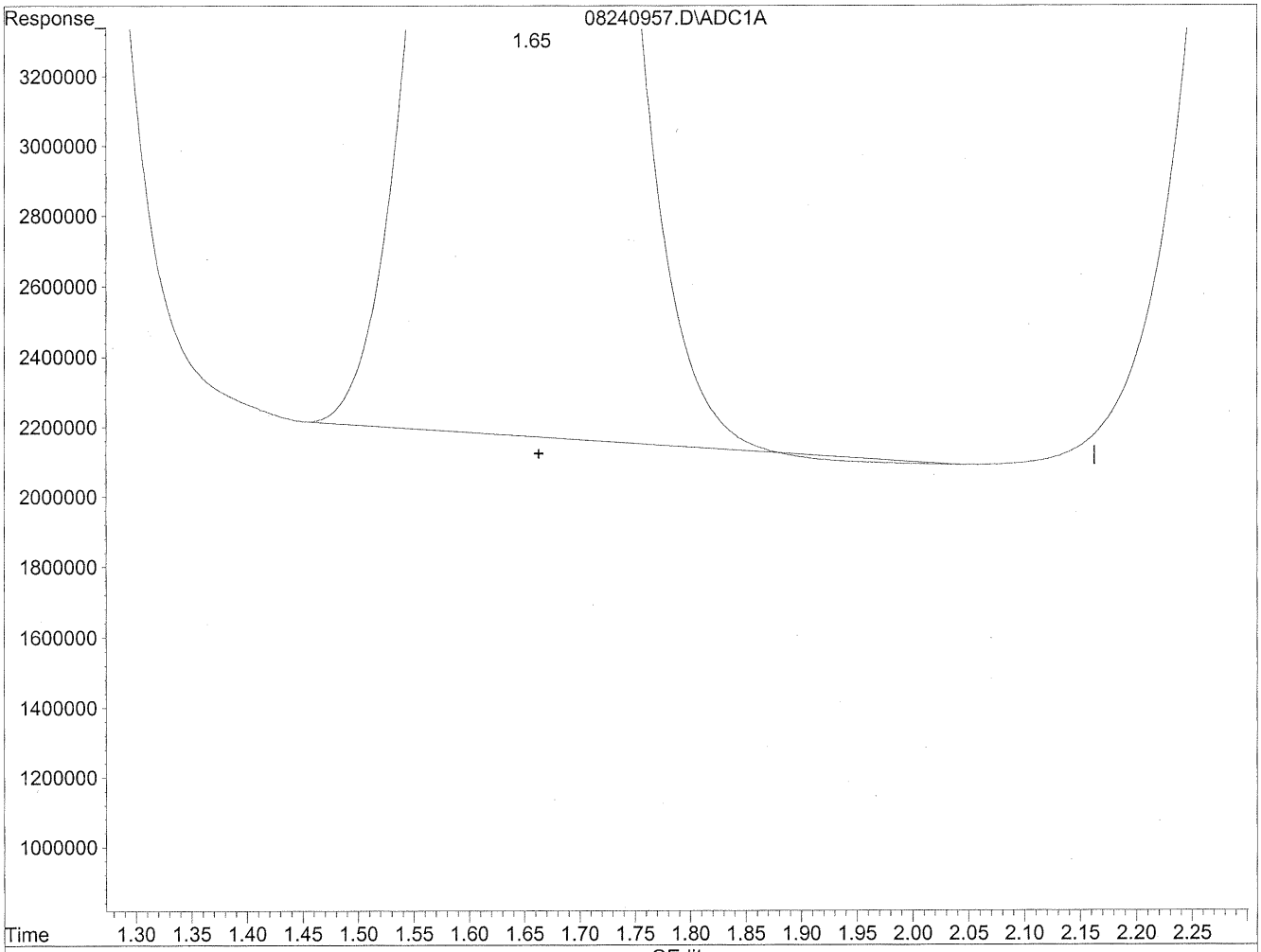
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.19	925252954	5040.015	ng/ml
2)	Acetaldehyde	1.65	634541585	4525.216	ng/mlm
3)	Propionaldehyde	2.97	50547819	473.759	ng/mlm
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	5.11	42508771	481.216	ng/mlm
6)	Benzaldehyde	6.56	38680814	587.236	ng/mlm
7)	Isovaleraldehyde	7.40	14921429	190.687	ng/mlm
8)	Valeraldehyde	7.87	96441628	1312.042	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	10.73	387432597	5753.058	ng/mlm
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/mld

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.65min 4510.756ng/ml

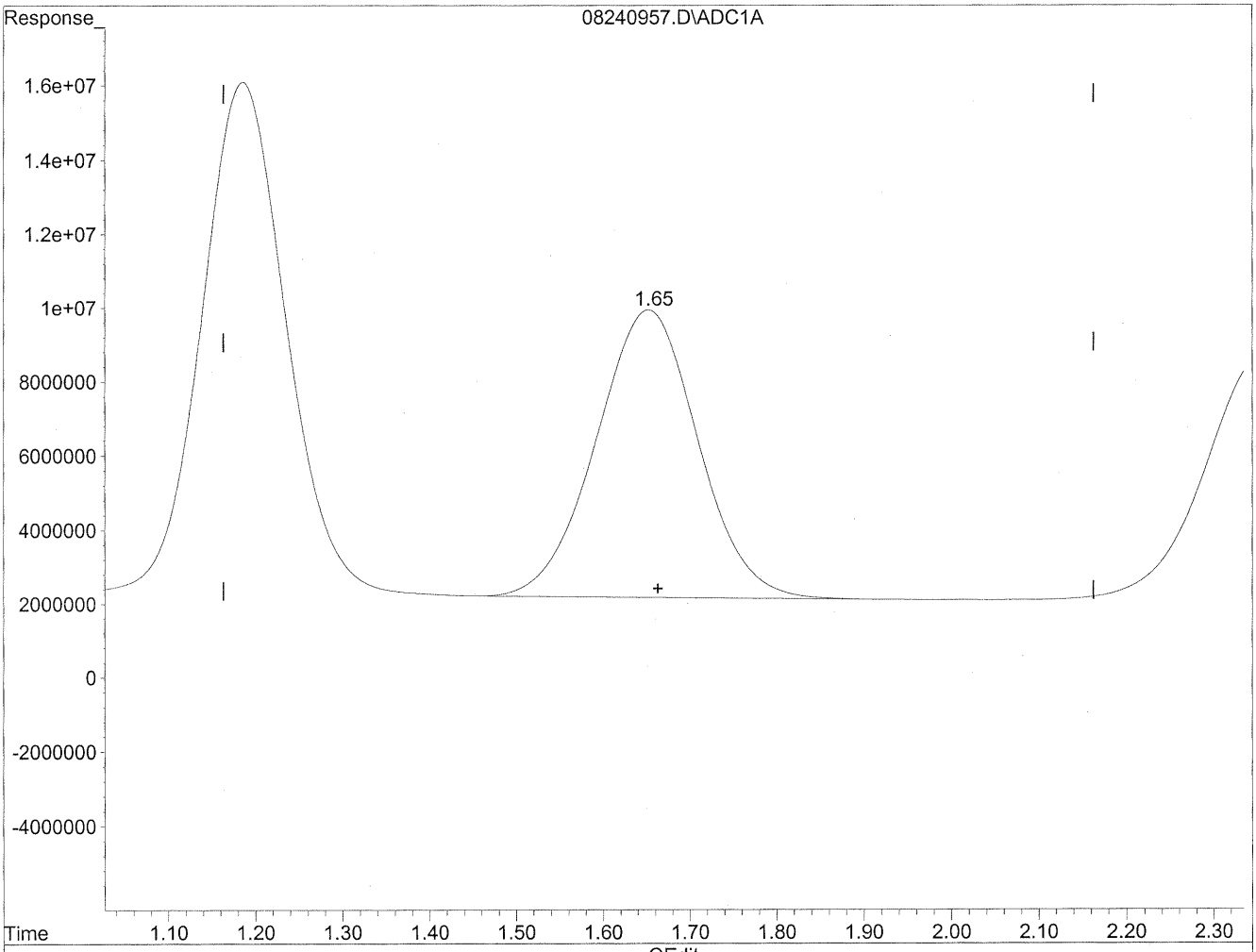
response 632514017

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.65min 4525.216ng/ml m

response 634541585

8/31/09
LC

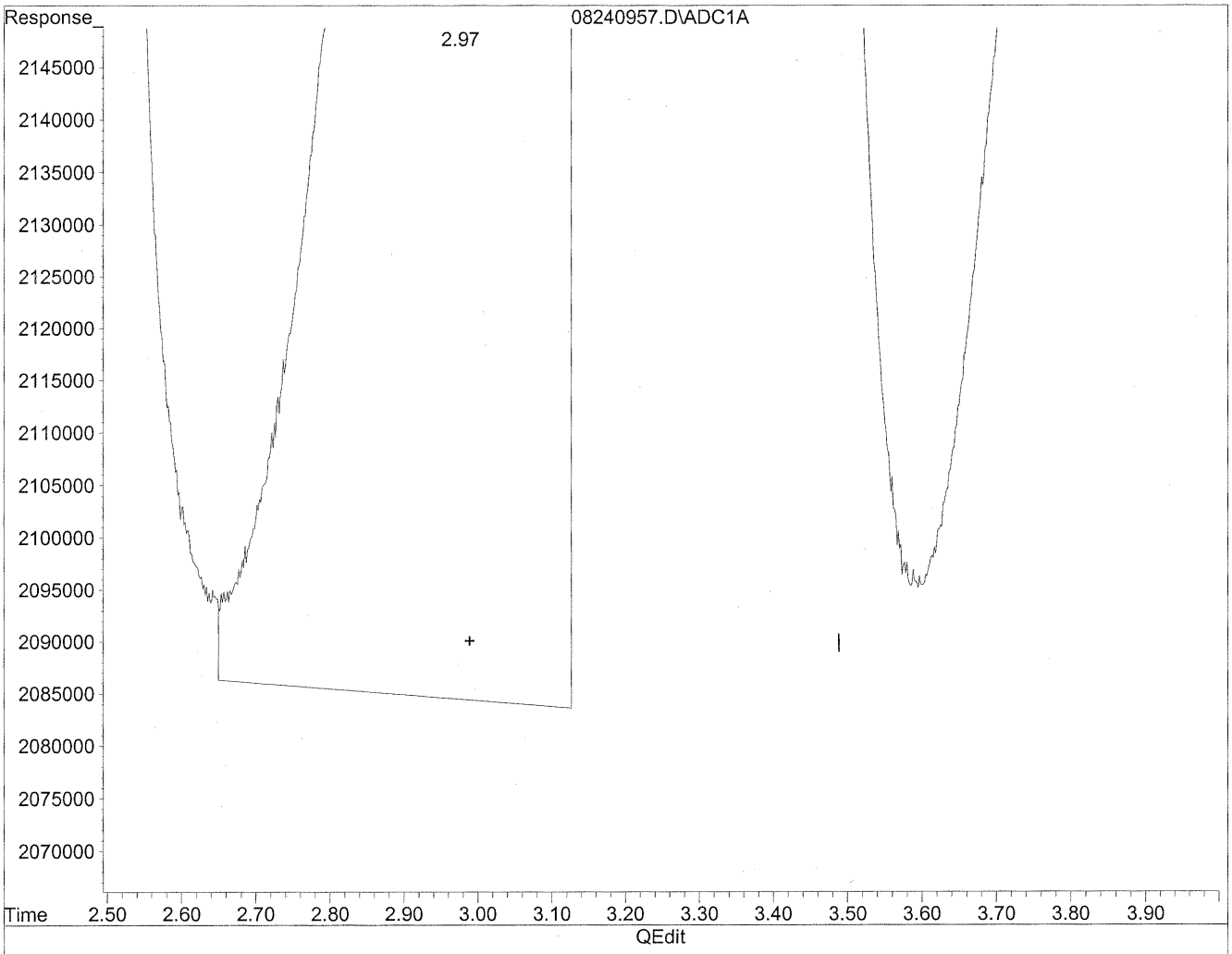
(+) = Expected Retention Time

08240957.D TO110709.M Sat Aug 29 14:55:51 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.97min 497.532ng/ml

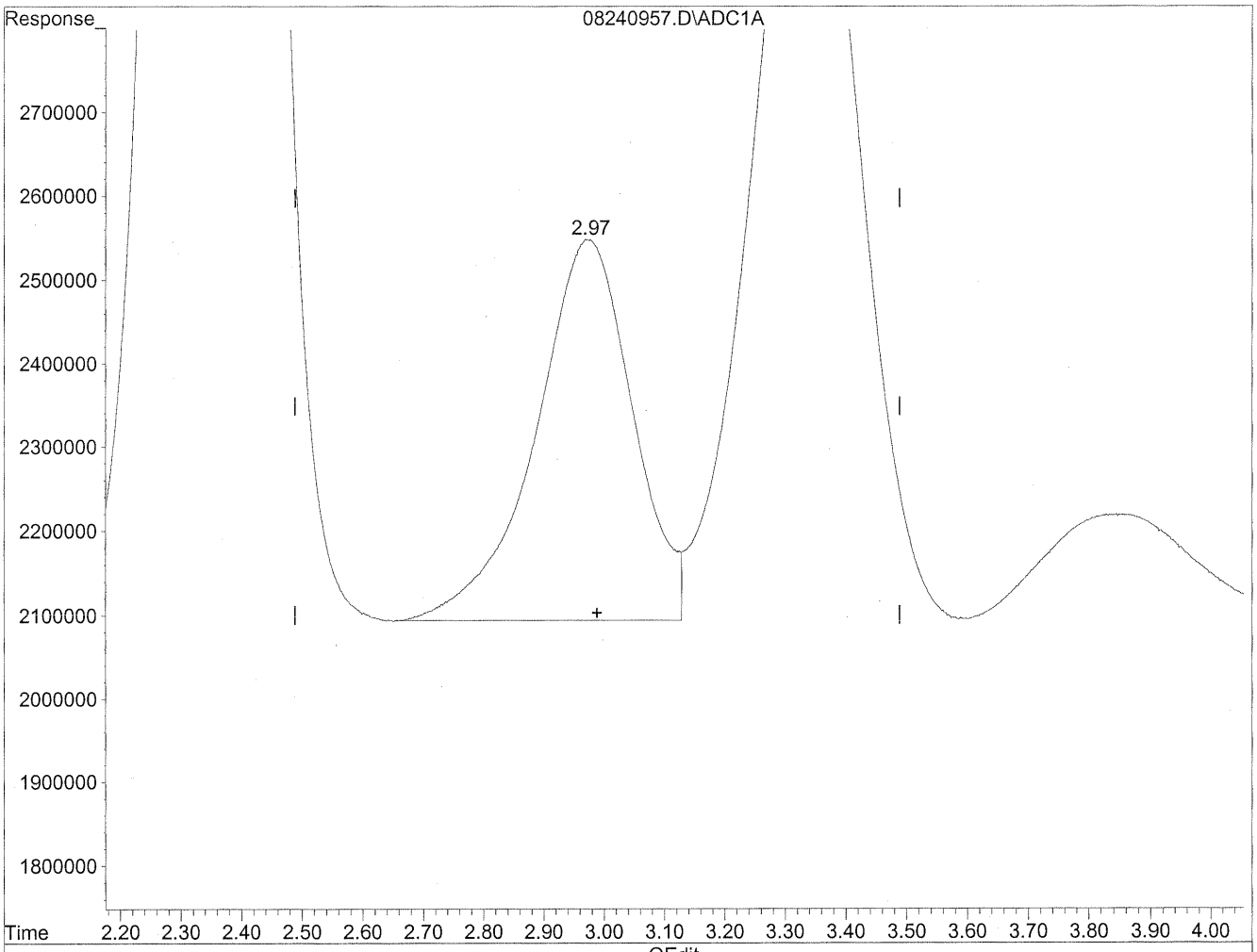
response 53084268

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.97min 473.759ng/ml m

response 50547819

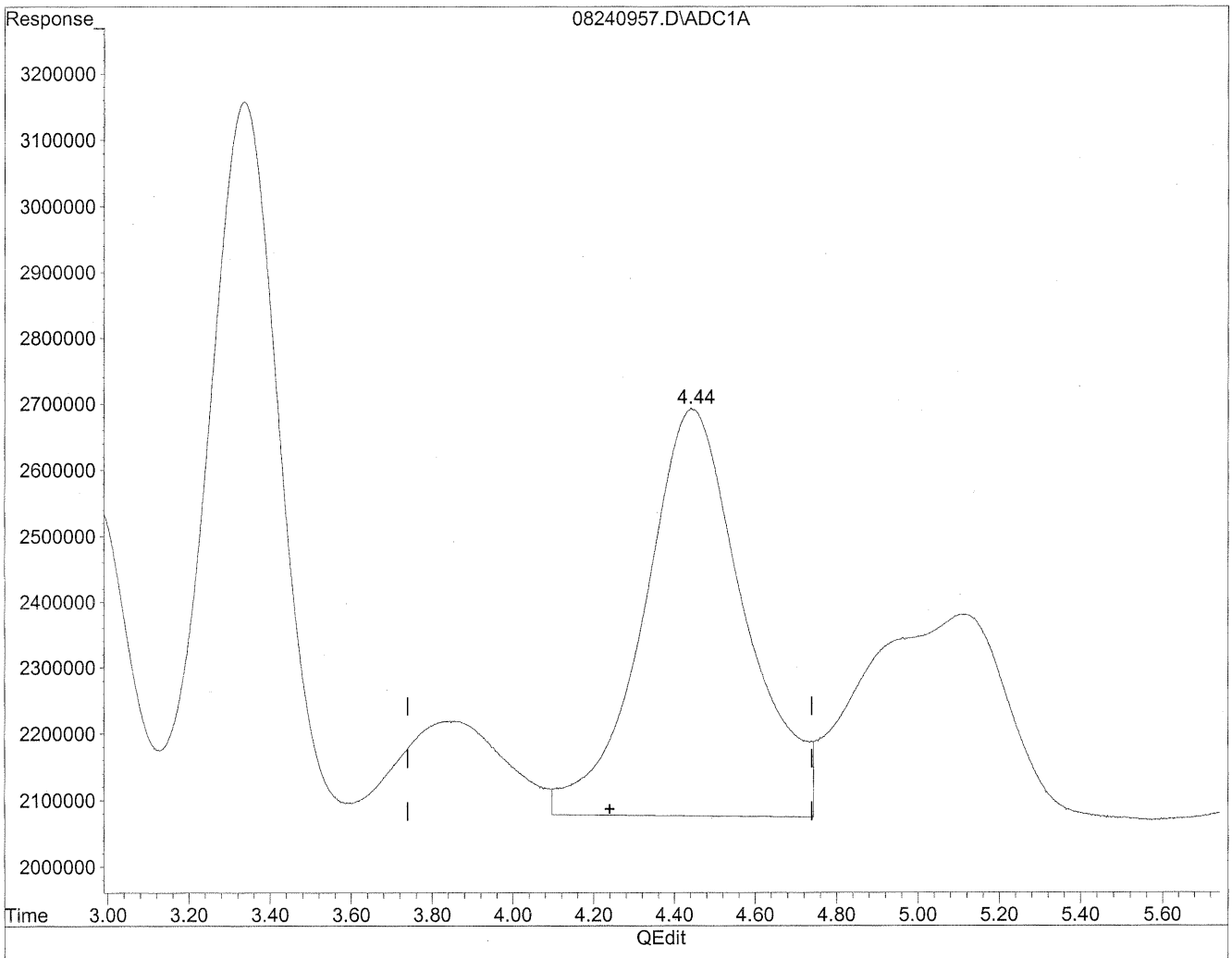
(+) = Expected Retention Time

08240957.D TO110709.M Sat Aug 29 14:56:07 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.44min 1092.848ng/ml

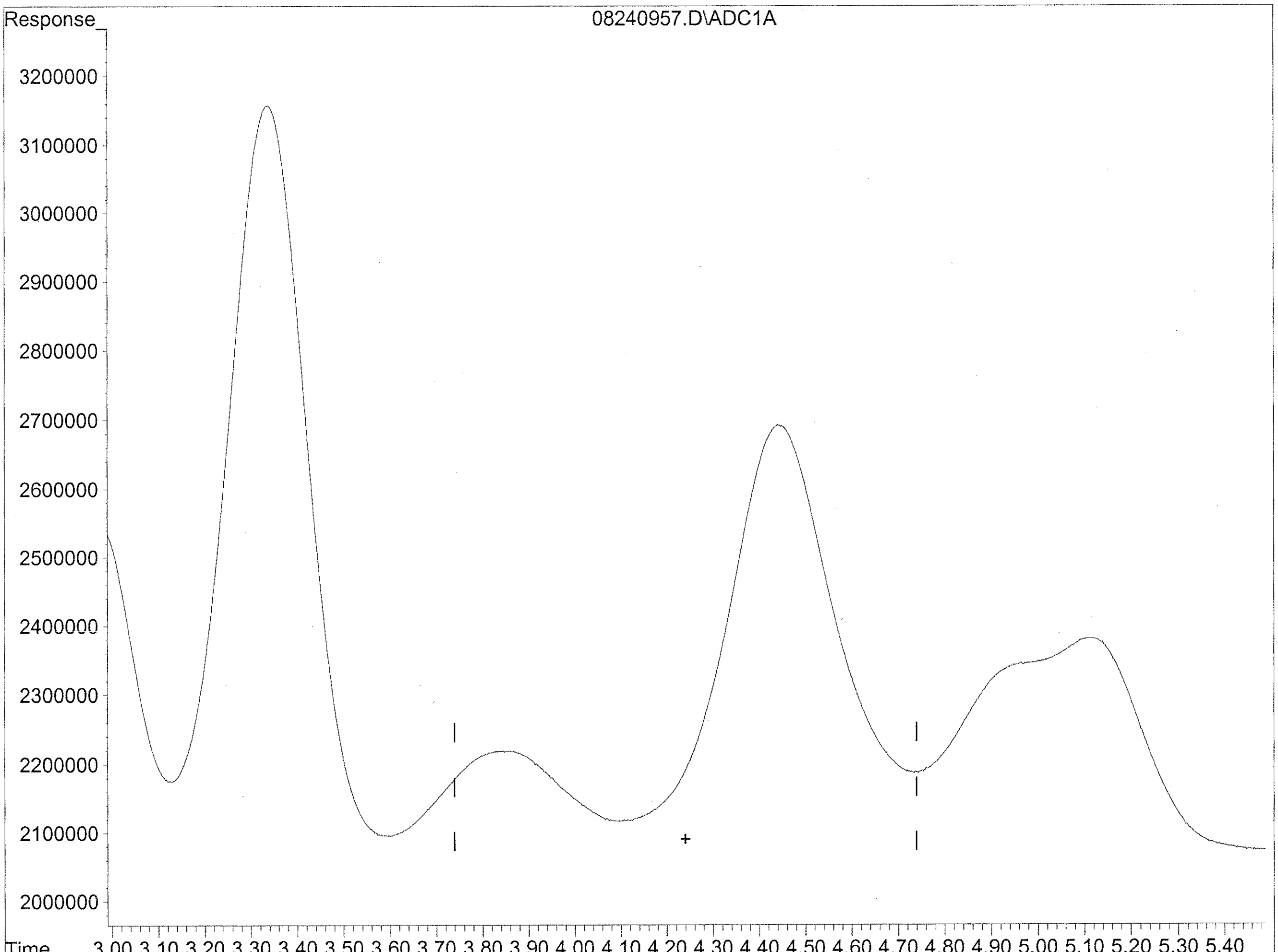
response 106460022

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



QEdit

(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

HC 8/29/09

MP

WB 8/31/09

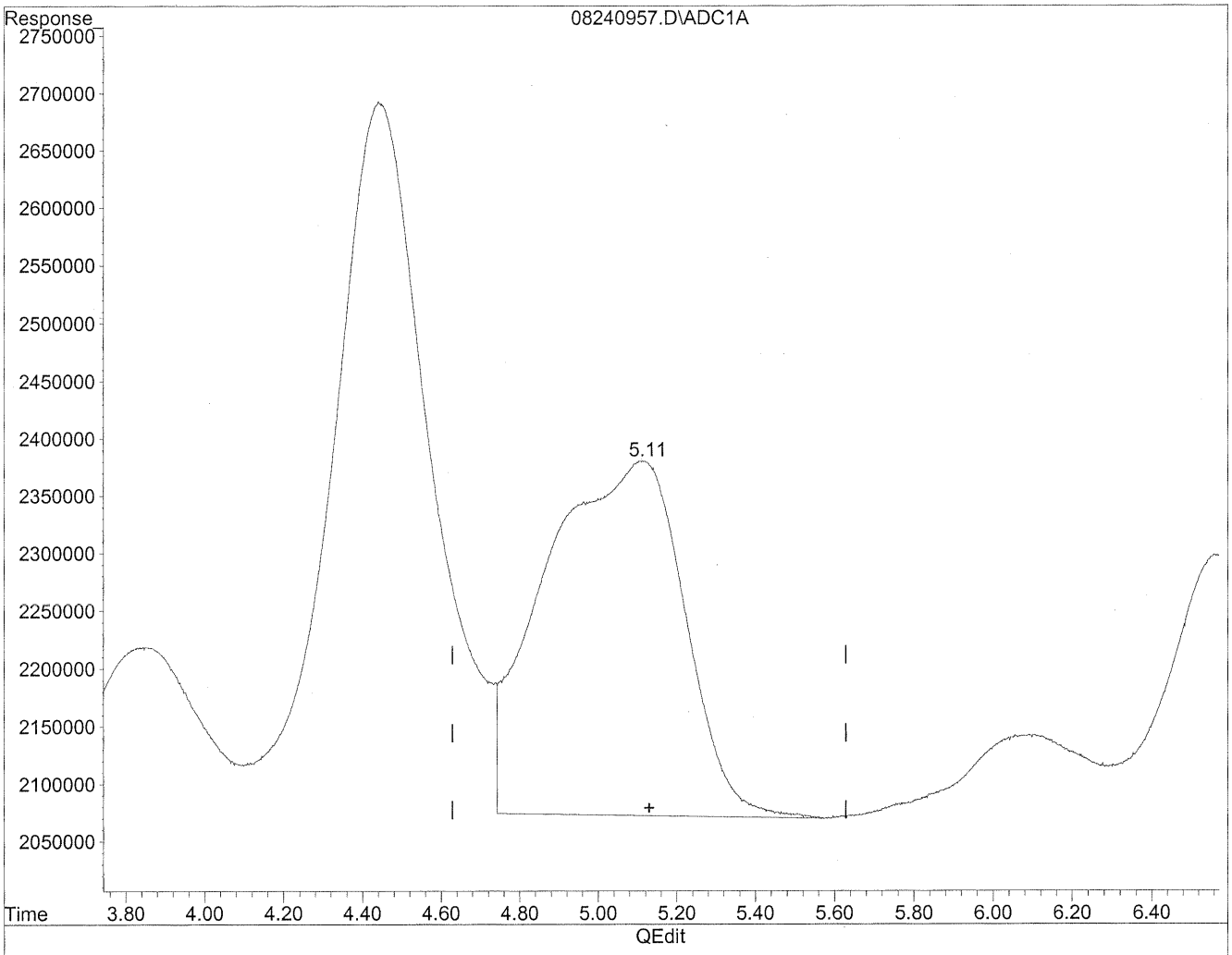
85

(+) = Expected Retention Time
08240957.D TO110709.M Sat Aug 29 14:56:13 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.11min 858.063ng/ml

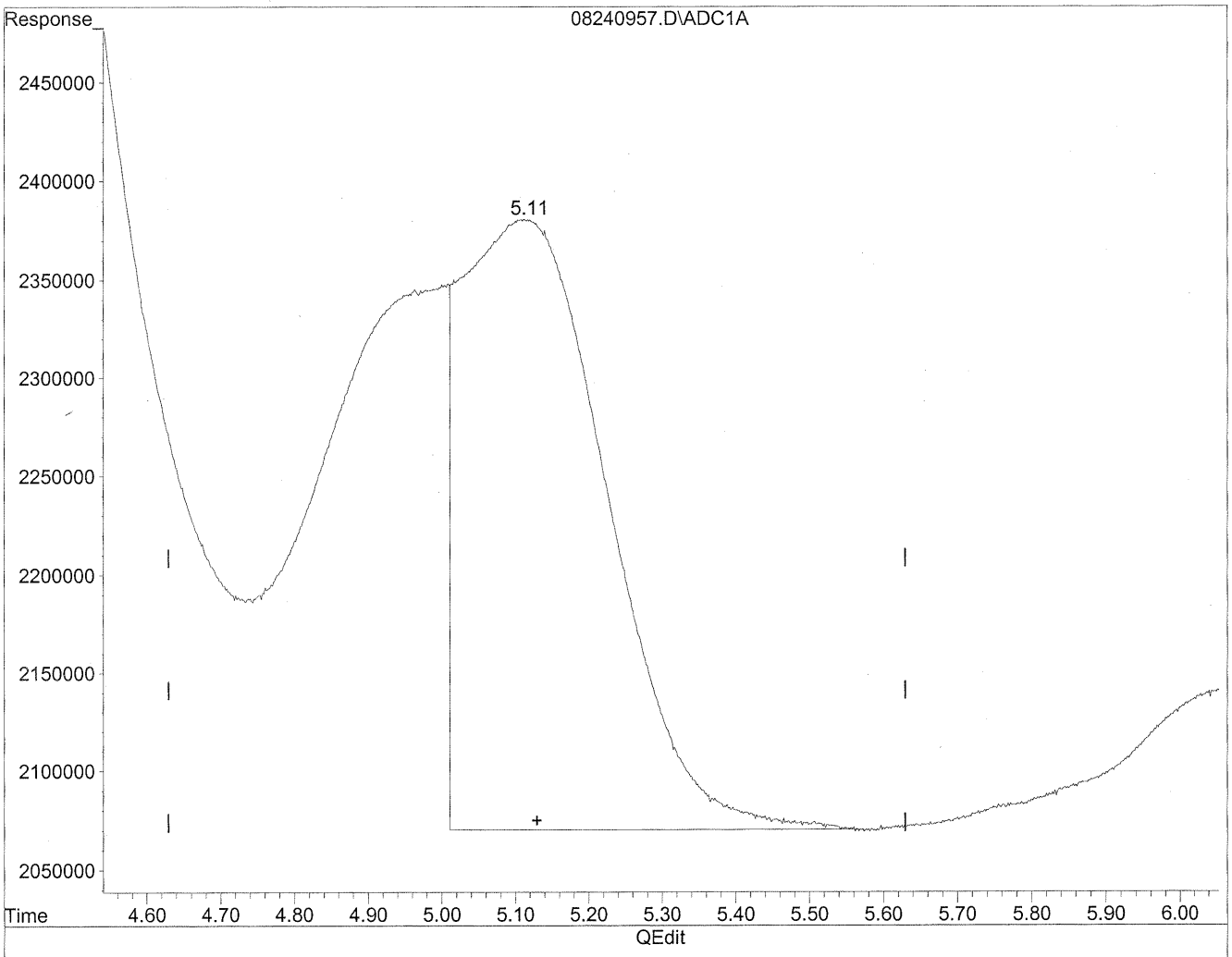
response 75797996

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.11min 481.216ng/ml m

response 42508771

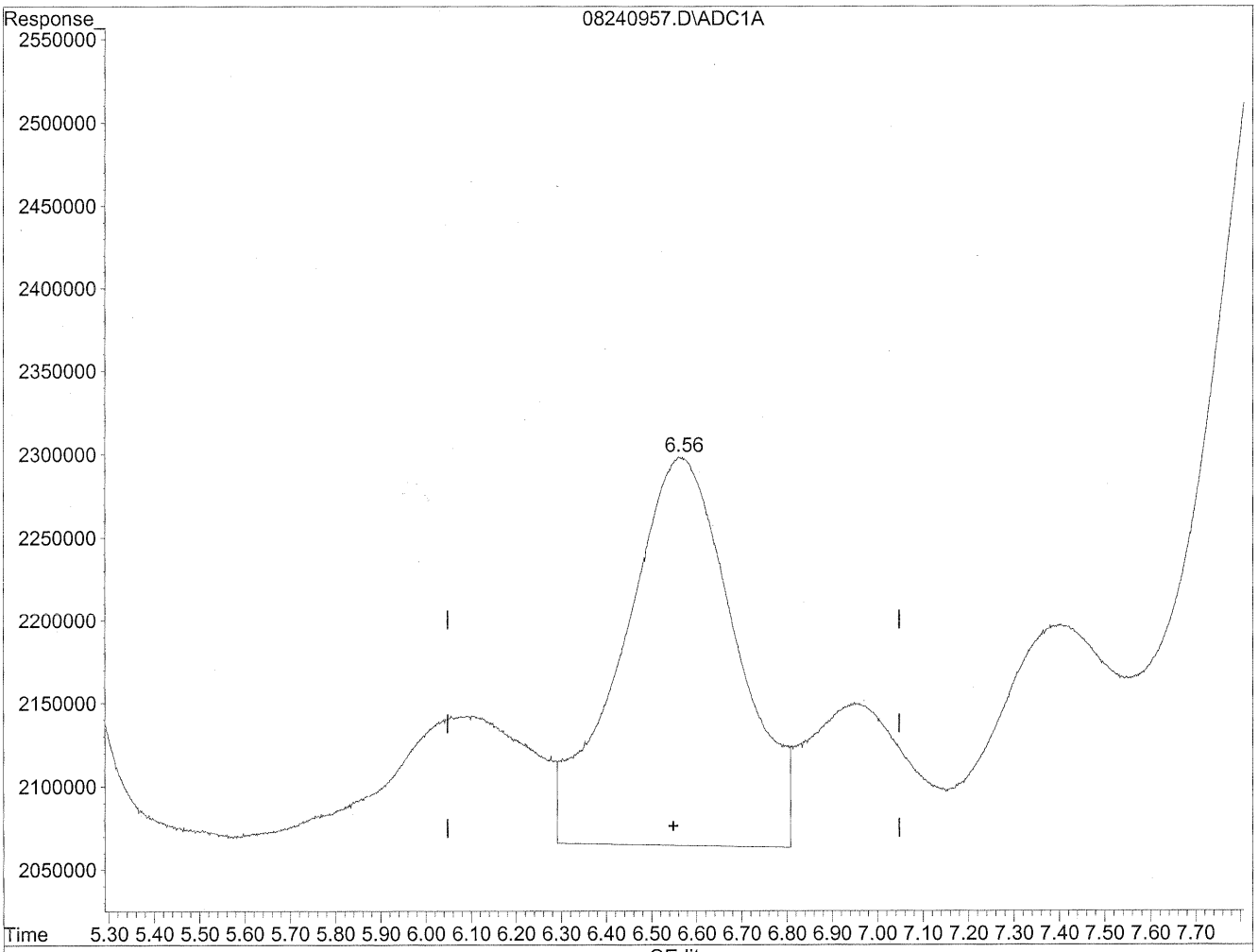
HC 8/29/09 SP 8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.57min 612.603ng/ml

response 40351752

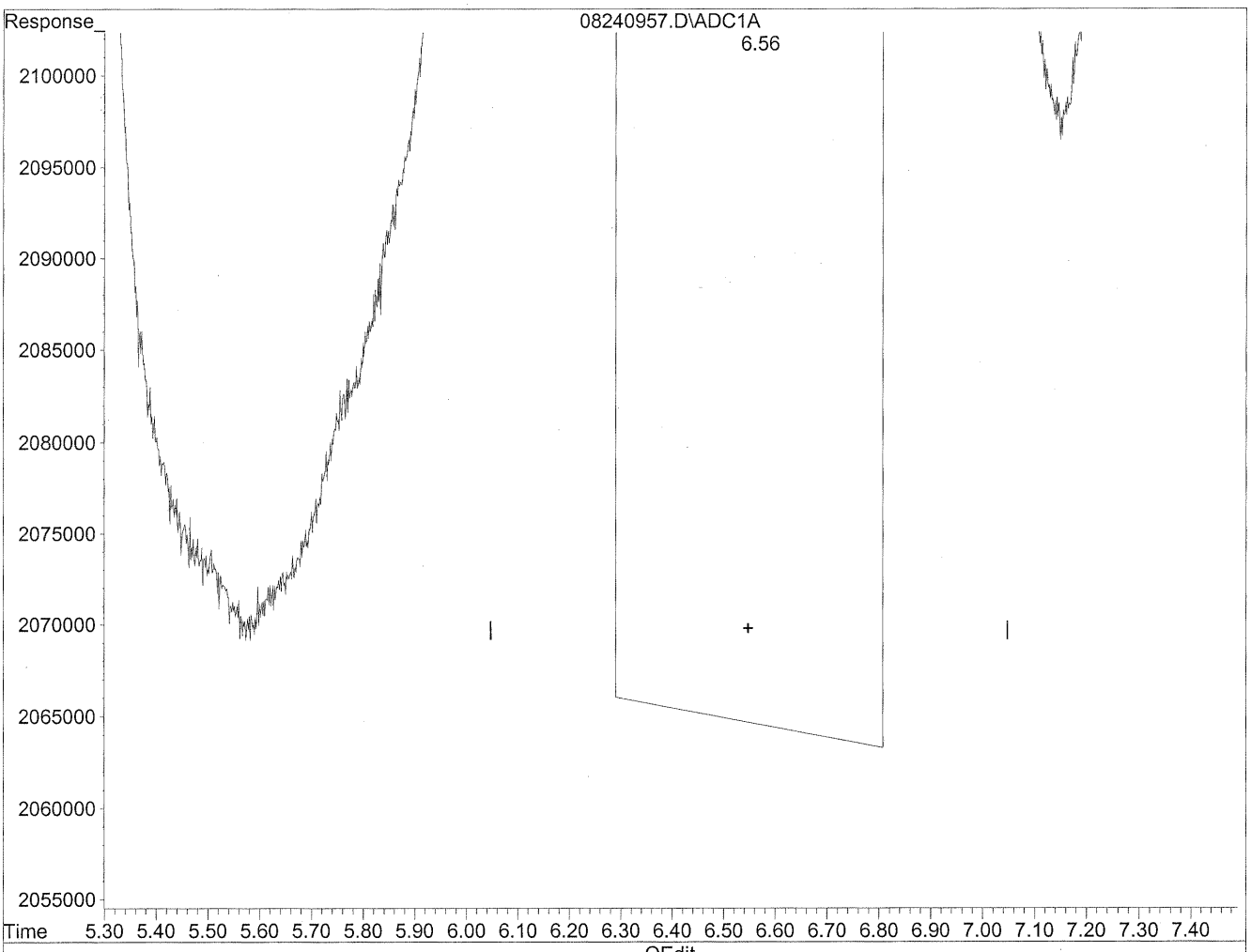
(+) = Expected Retention Time

08240957.D TO110709.M Sat Aug 29 14:56:47 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.57min 612.603ng/ml

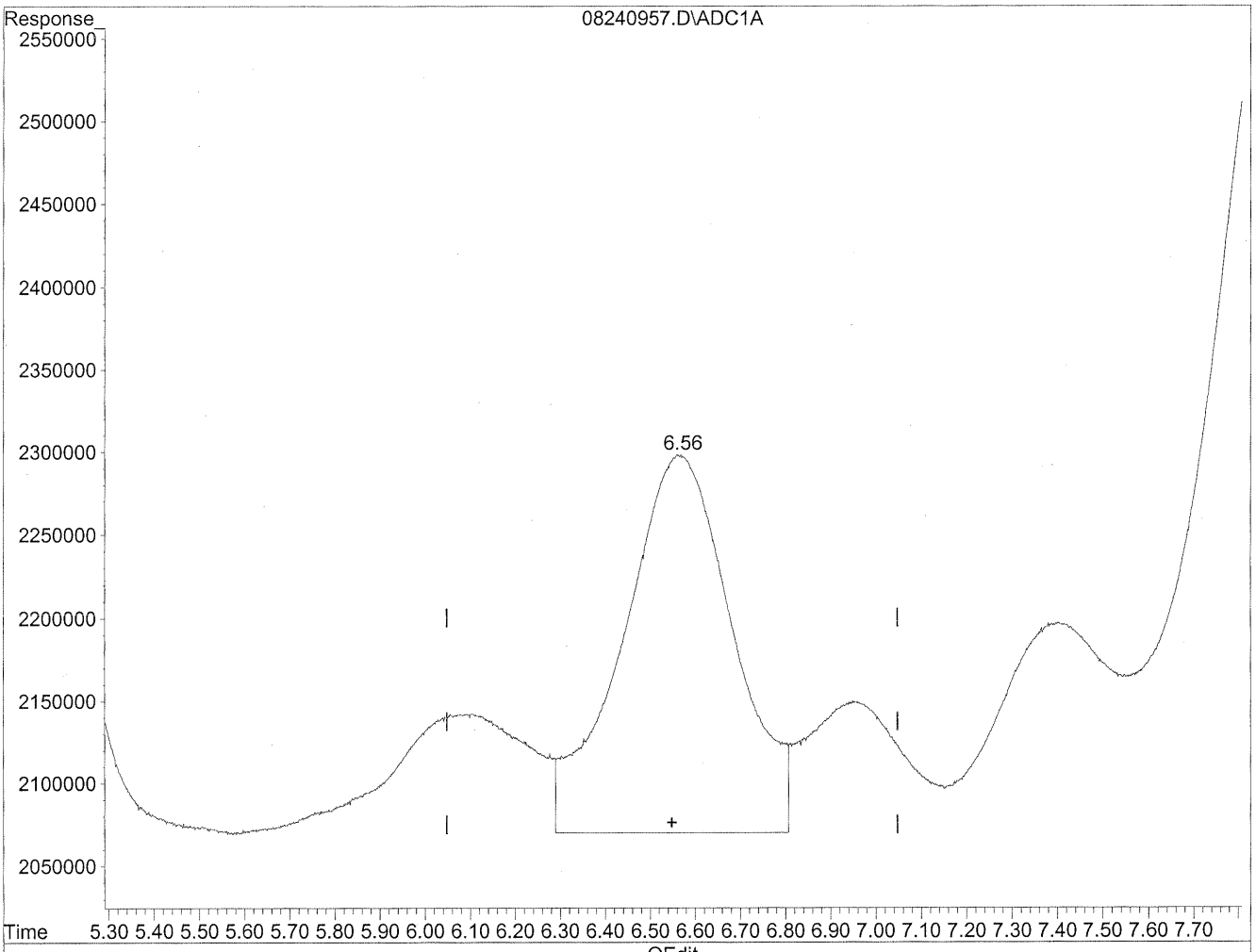
response 40351752

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.56min 587.236ng/ml m

response 38680814

HC 8/29/09 BC

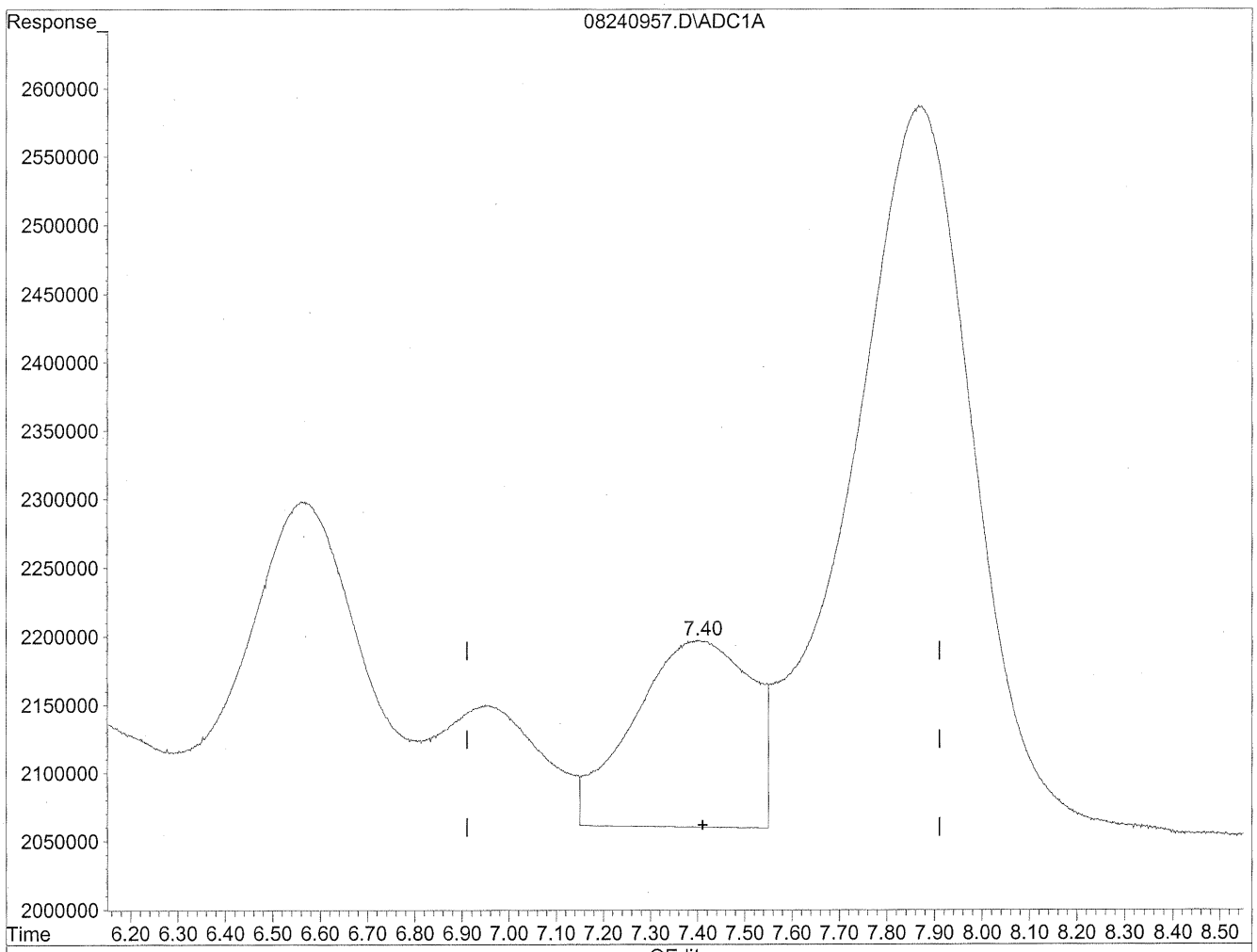
LM 8/31/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.40min 304.365ng/ml

response 23816862

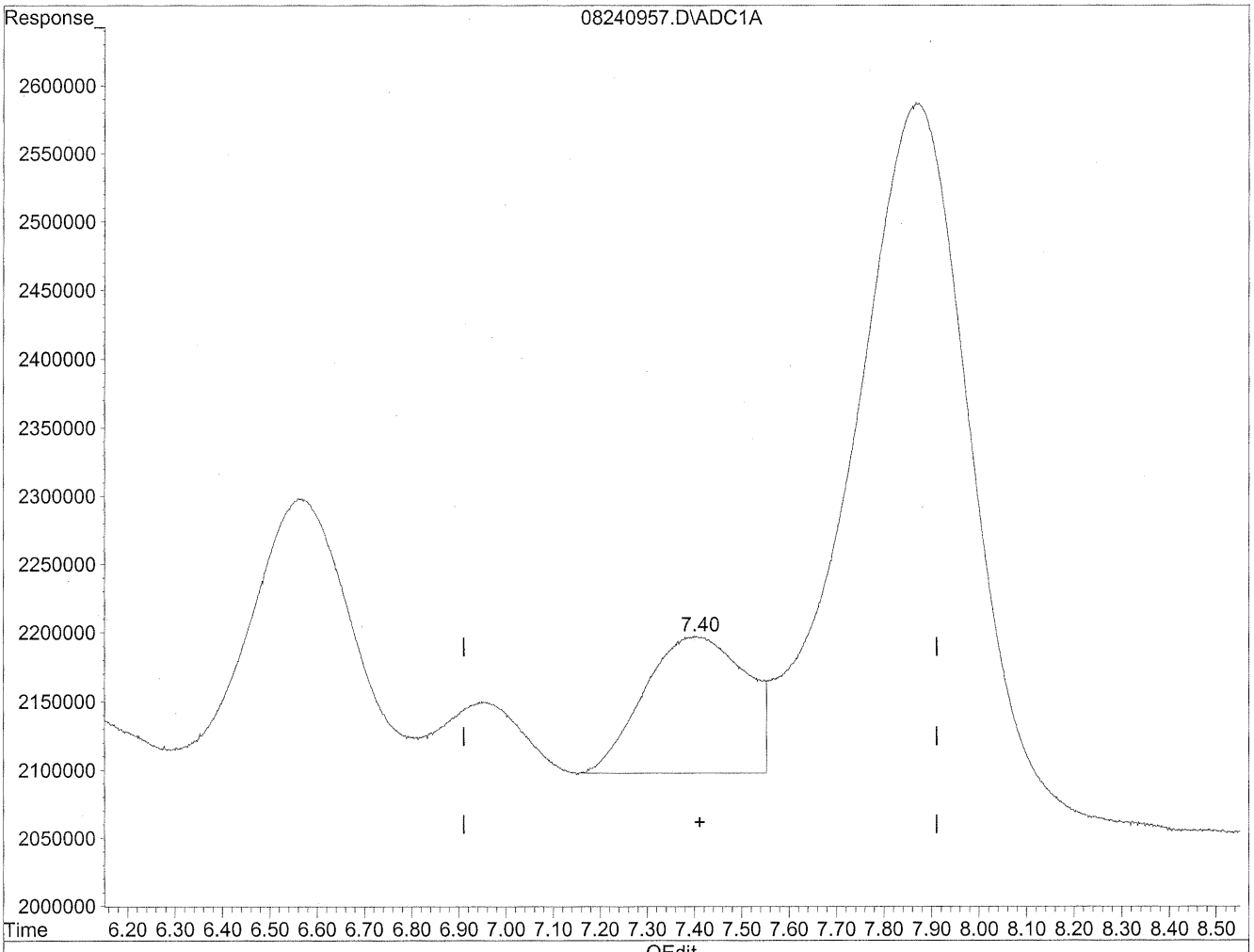
(+) = Expected Retention Time

08240957.D TO110709.M Sat Aug 29 14:57:17 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.40min 190.687ng/ml m

response 14921429

HC 8/29/09 BC

CW 9/3/09

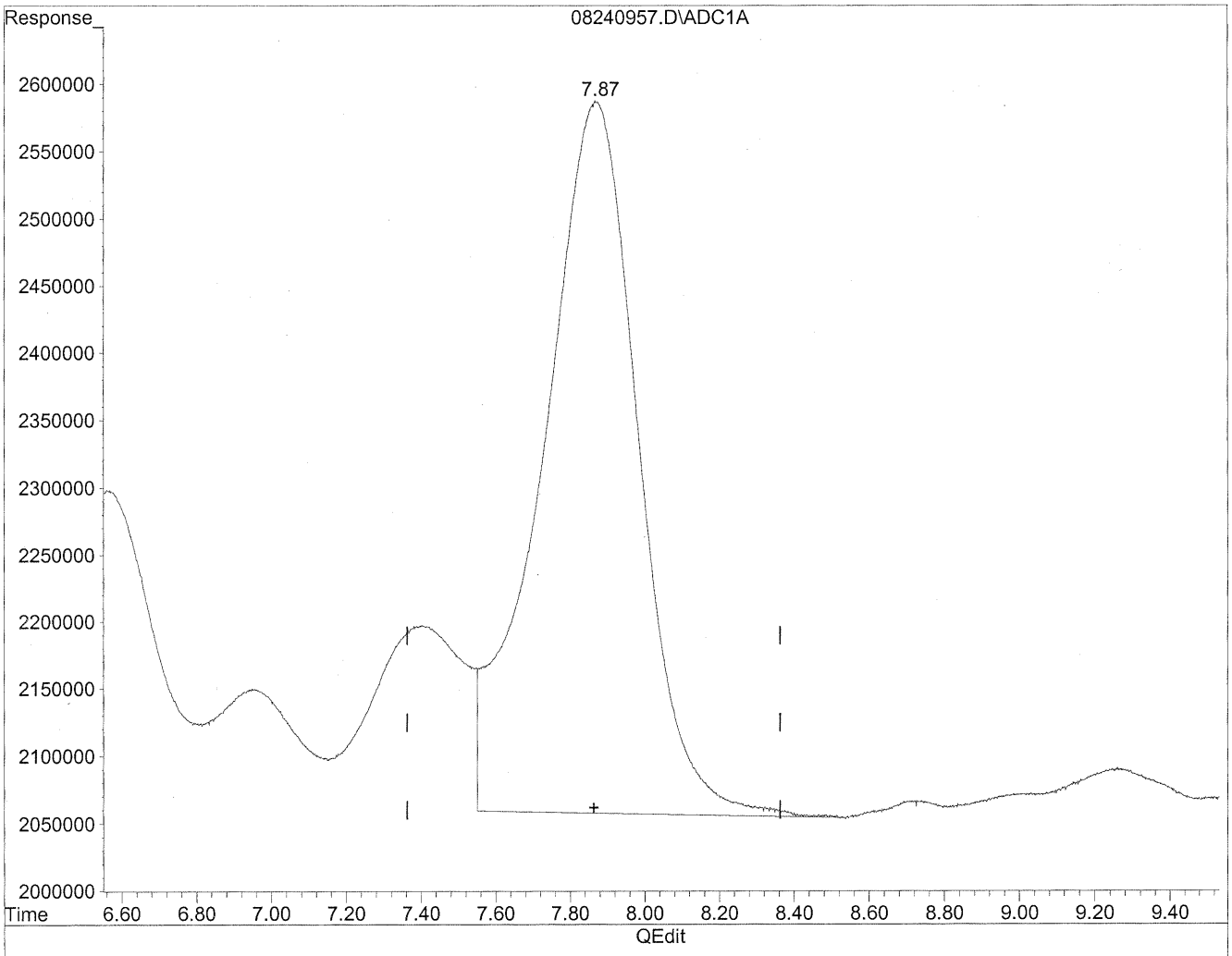
92

(+) = Expected Retention Time
08240957.D TO110709.M Sat Aug 29 14:57:29 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.87min 1291.254ng/ml

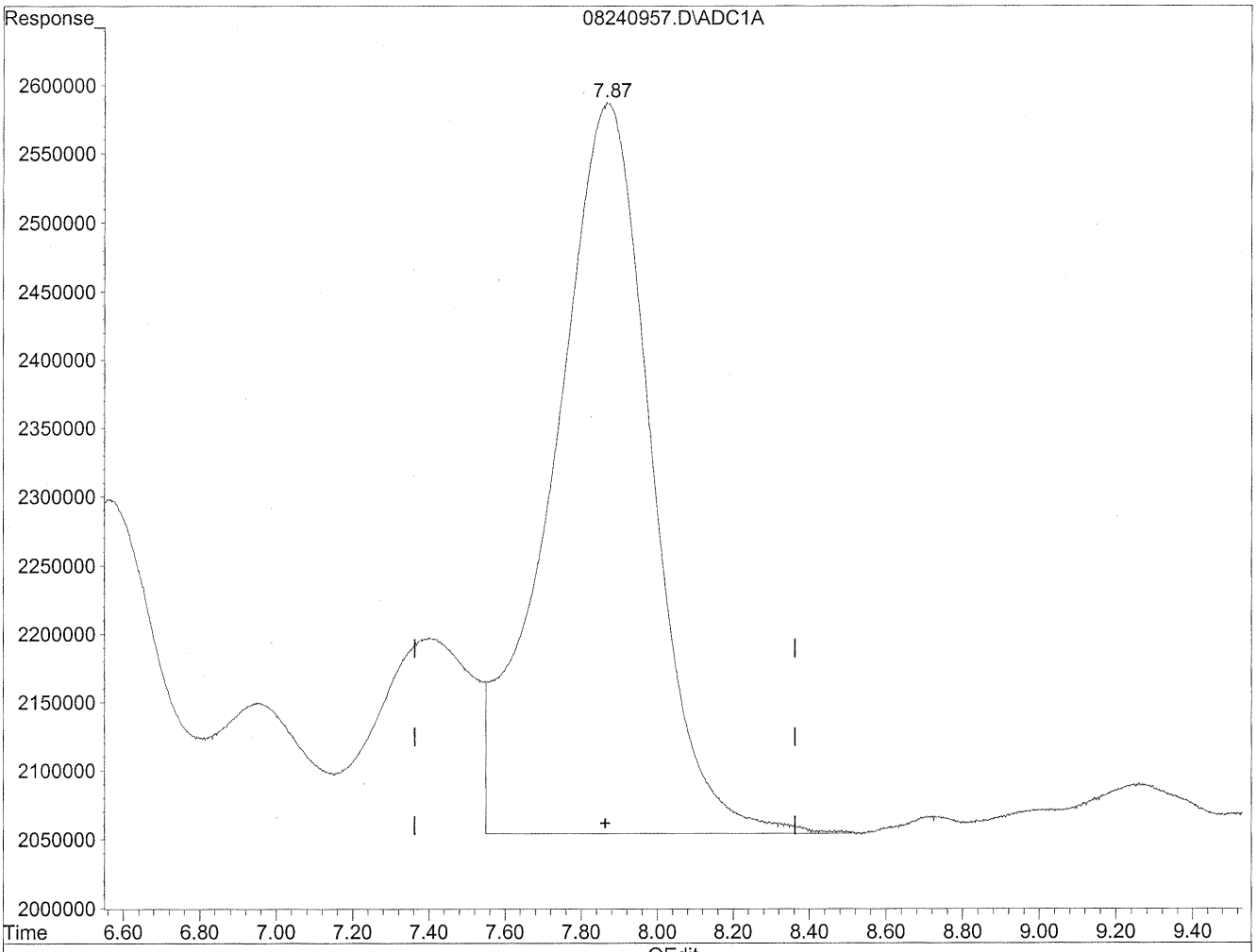
response 94913663

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.87min 1312.042ng/ml m

response 96441628

8/29/09 LC

8/31/09

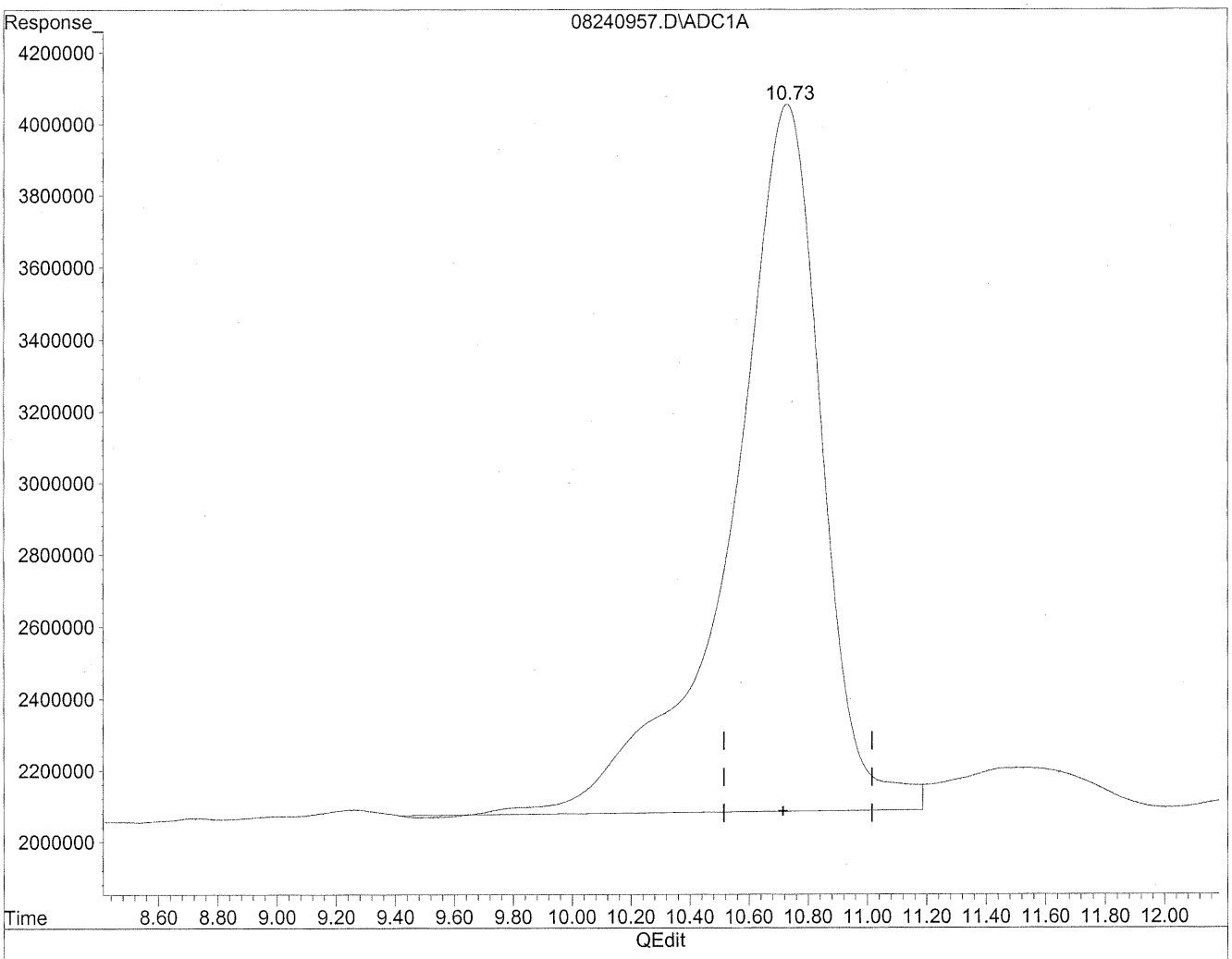
(+) = Expected Retention Time

08240957.D TO110709.M Sat Aug 29 14:57:50 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

10.73min 6273.632ng/ml

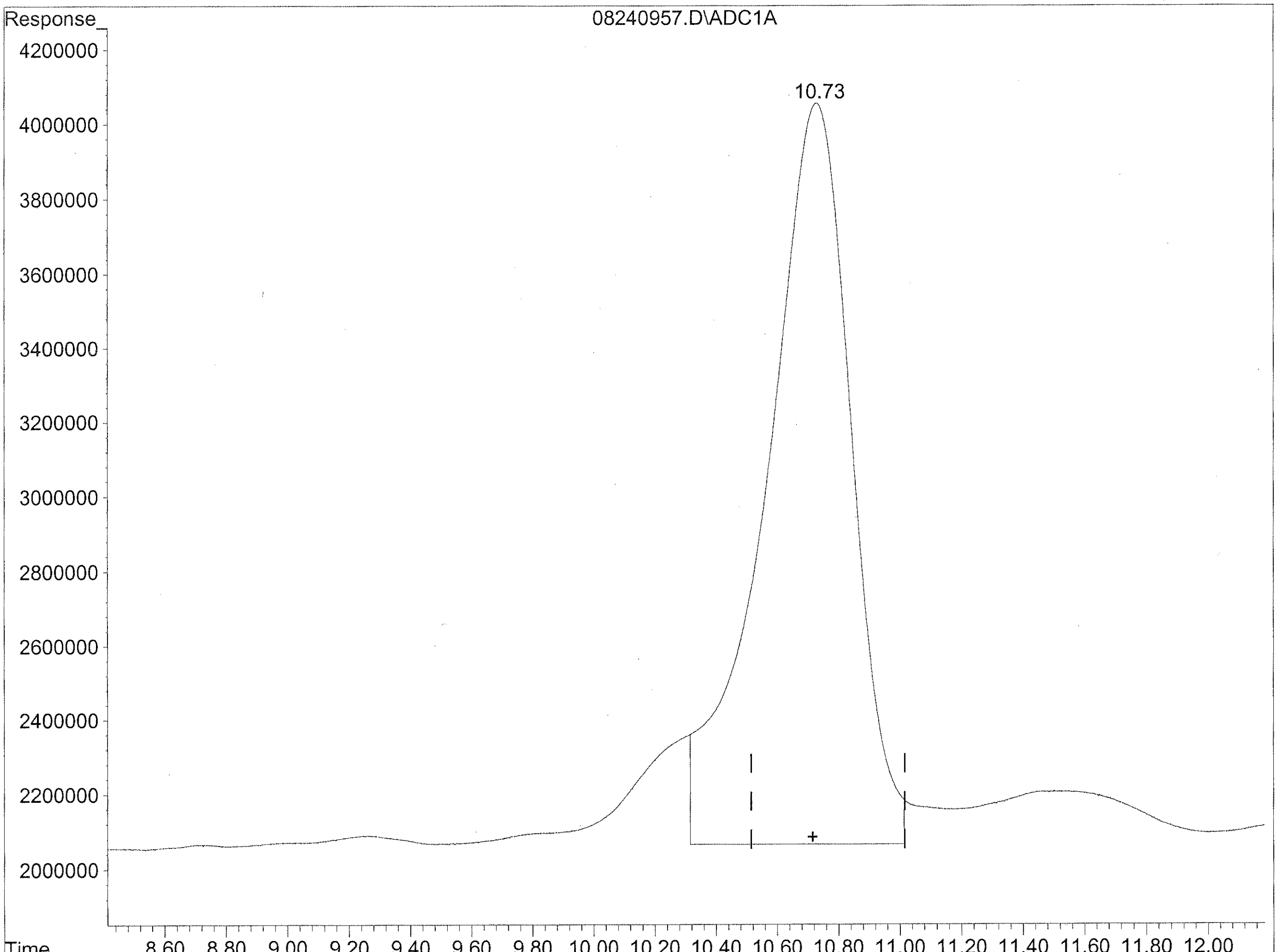
response 422489980

(+) = Expected Retention Time

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D       Vial: 54
Acq On    : 25 Aug 2009   2:32 am                       Operator: HC
Sample    : P0902912-004 front 1.0ml                    Inst    : LC 01
Misc      :                                             Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Aug 25  7:20 19109  Quant Results File: TO110709.RES

Method        : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title         : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update   : Sat Aug 29 12:41:27 2009
Response via  : Multiple Level Calibration
```



QEdit

(11) Hexaldehyde

10.73min 5753.058ng/ml m

response 387432597

HC 8/29/09 5H,BC

WA 8/31/09

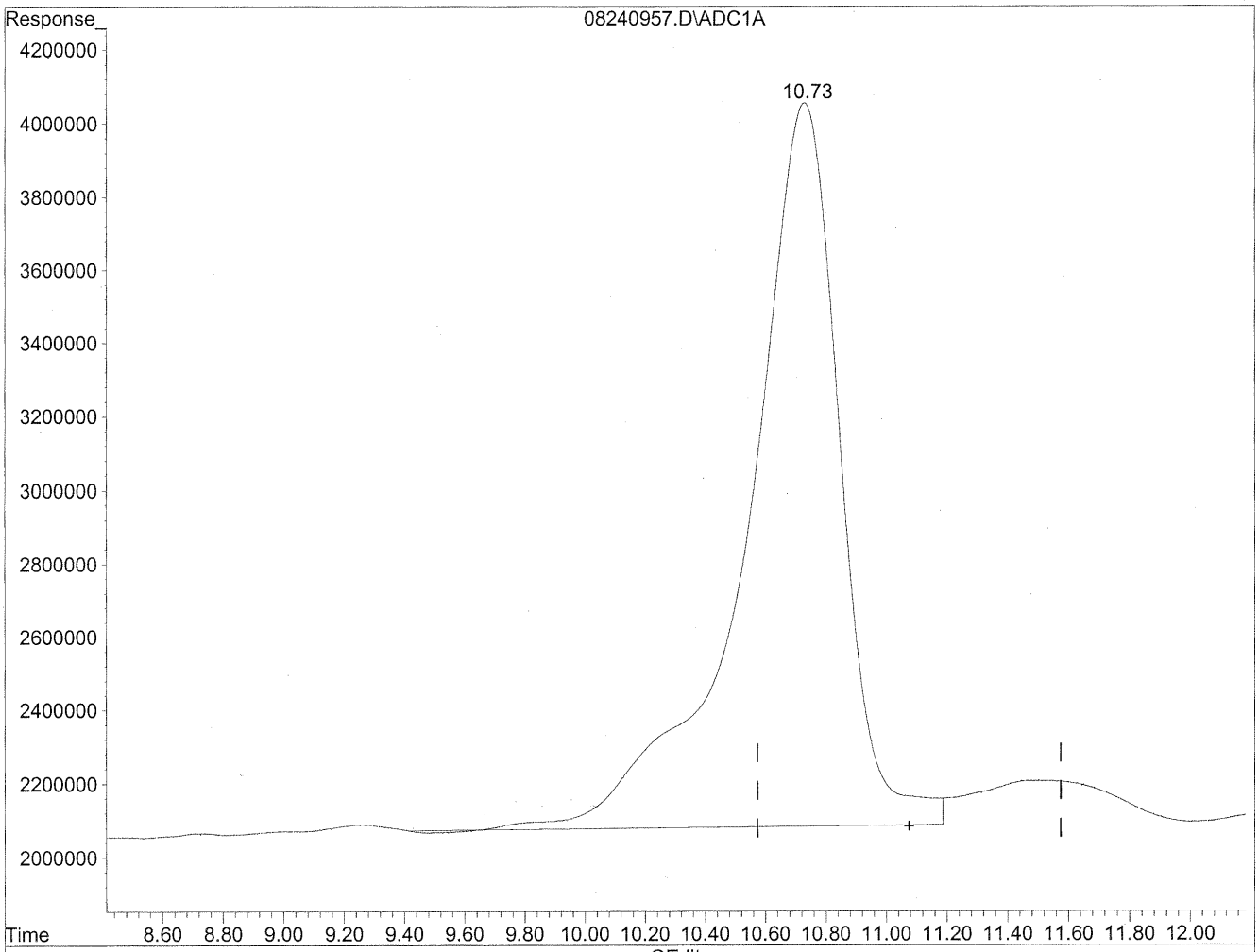
(+) = Expected Retention Time

08240957.D TO110709.M Sat Aug 29 14:58:37 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

10.73min 8619.889ng/ml

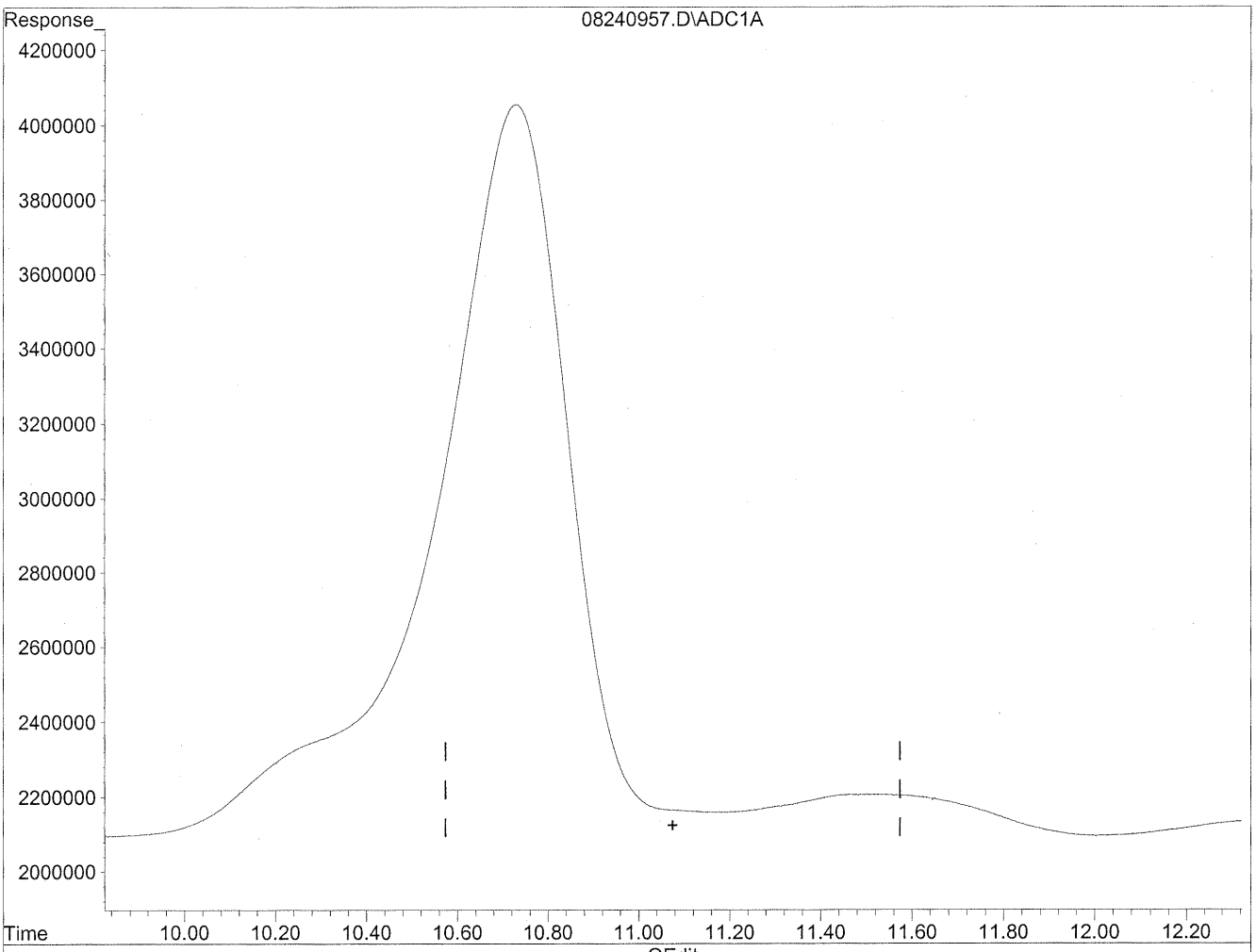
response 422489980

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240957.D Vial: 54
Acq On : 25 Aug 2009 2:32 am Operator: HC
Sample : P0902912-004 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 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

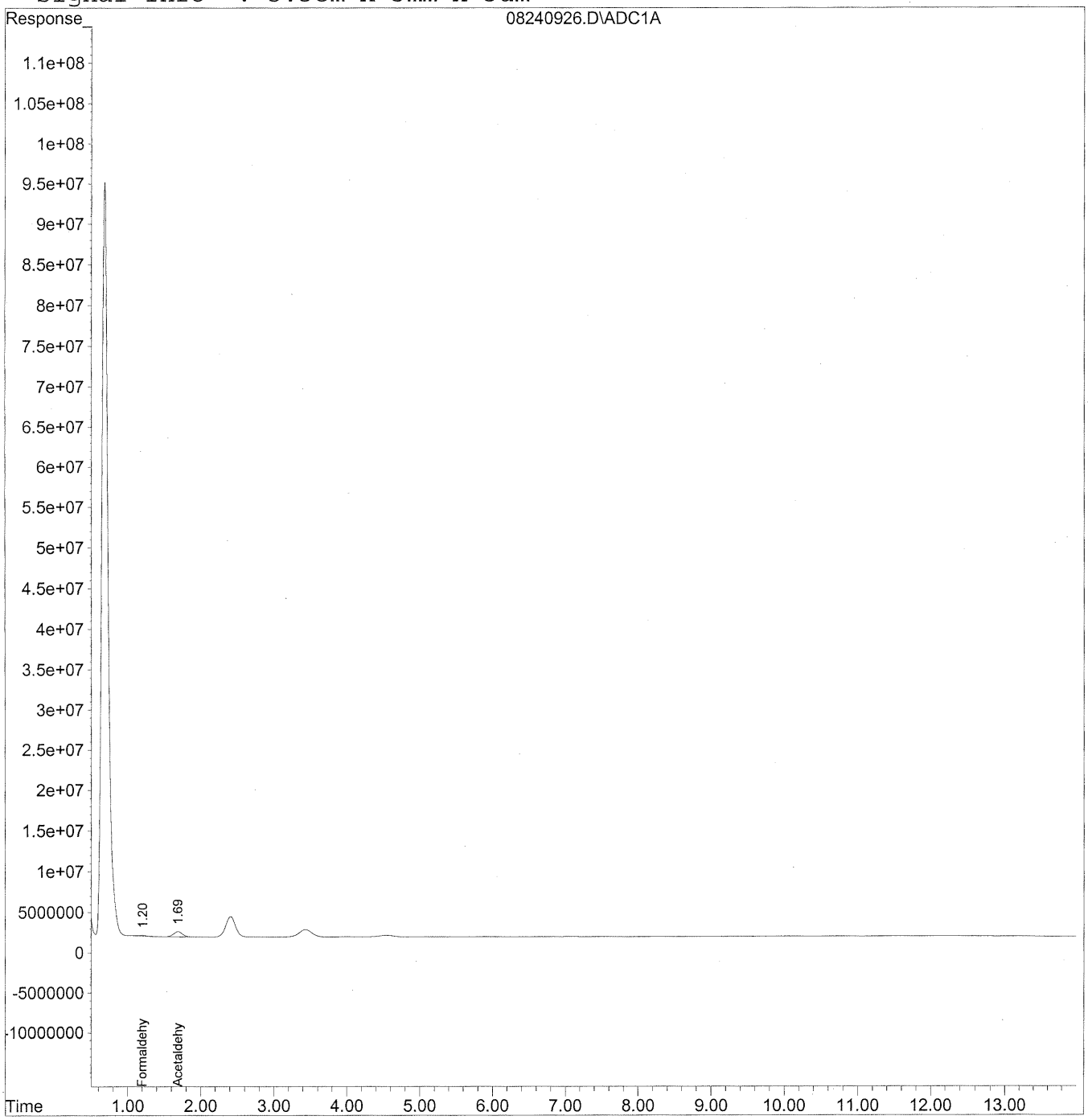
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240926.D Vial: 24
Acq On : 24 Aug 2009 6:46 pm Operator: HC
Sample : P0902912-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 13:37 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 25 10:21:57 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\24\08240926.D Vial: 24
Acq On : 24 Aug 2009 6:46 pm Operator: HC
Sample : P0902912-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 13:37 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Aug 25 10:21:57 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

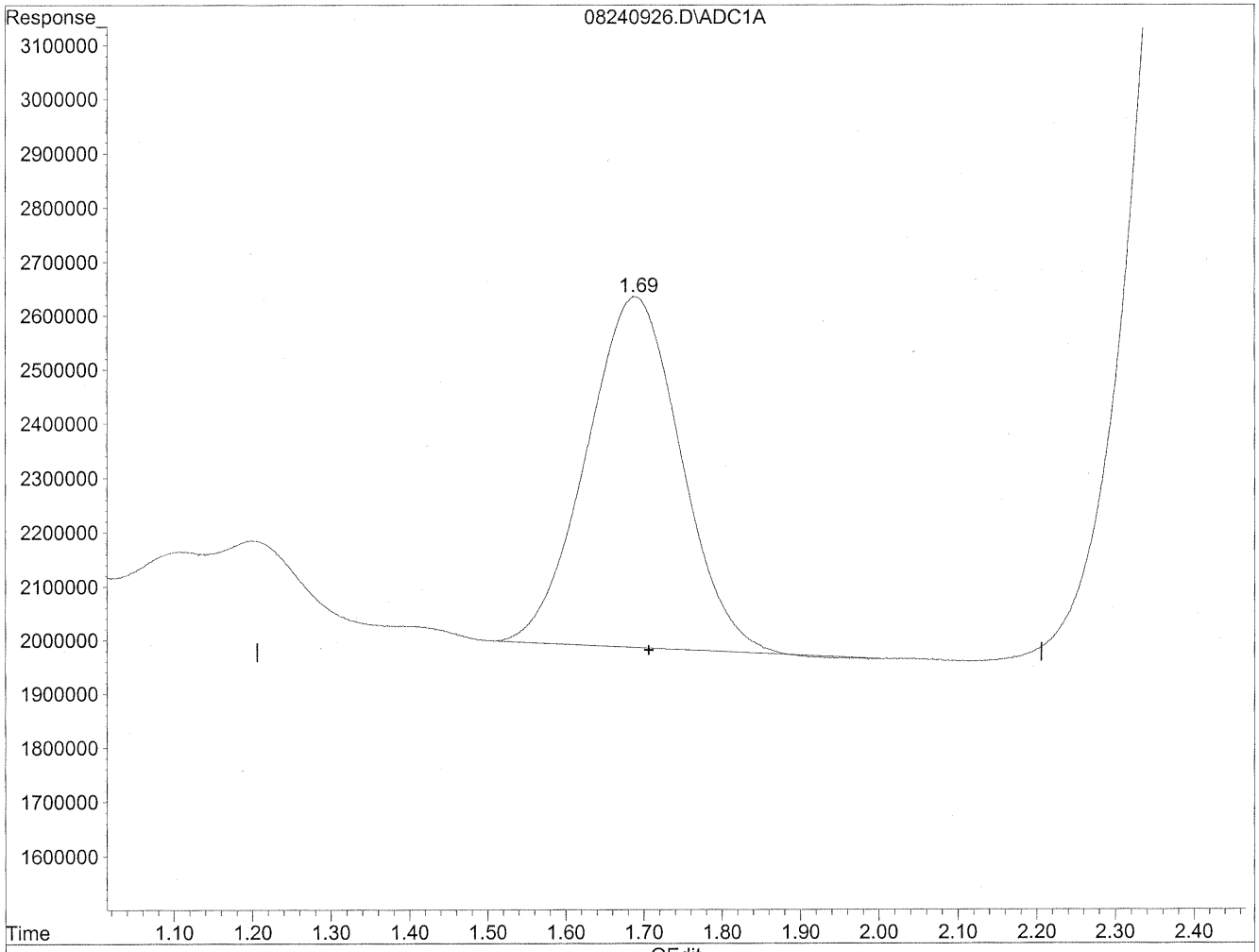
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.20	10221777	55.680	ng/ml
2)	Acetaldehyde	1.69	54788122	390.720	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/mld
4)	Crotonaldehyde	0.00	0	N.D.	ng/mld
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	0	N.D.	ng/ml
8)	Valeraldehyde	0.00	0	N.D.	ng/ml
9)	o-Tolualdehyde	0.00	0	N.D.	ng/ml
10)	m,p-Tolualdehyde	0.00	0	N.D.	ng/ml
11)	Hexaldehyde	0.00	0	N.D.	ng/ml
12)	2,5-Dimethylbenzaldehyde	0.00	0	N.D.	ng/ml

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240926.D Vial: 24
Acq On : 24 Aug 2009 6:46 pm Operator: HC
Sample : P0902912-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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

1.69min 387.989ng/ml

response 54405166

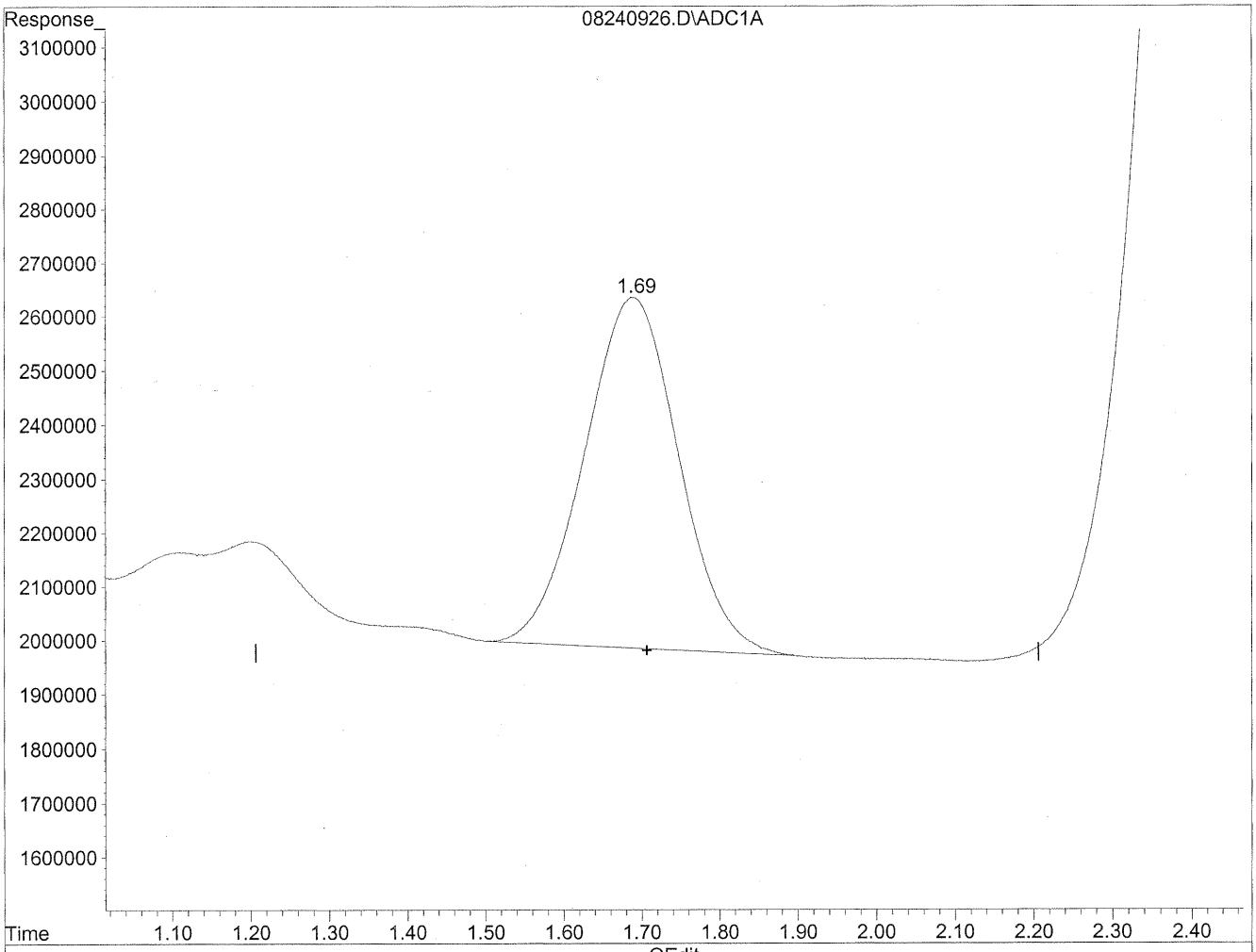
101

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240926.D Vial: 24
Acq On : 24 Aug 2009 6:46 pm Operator: HC
Sample : P0902912-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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

1.69min 390.720ng/ml m

response 54788122

HC 8/29/09 1C

8/31/09

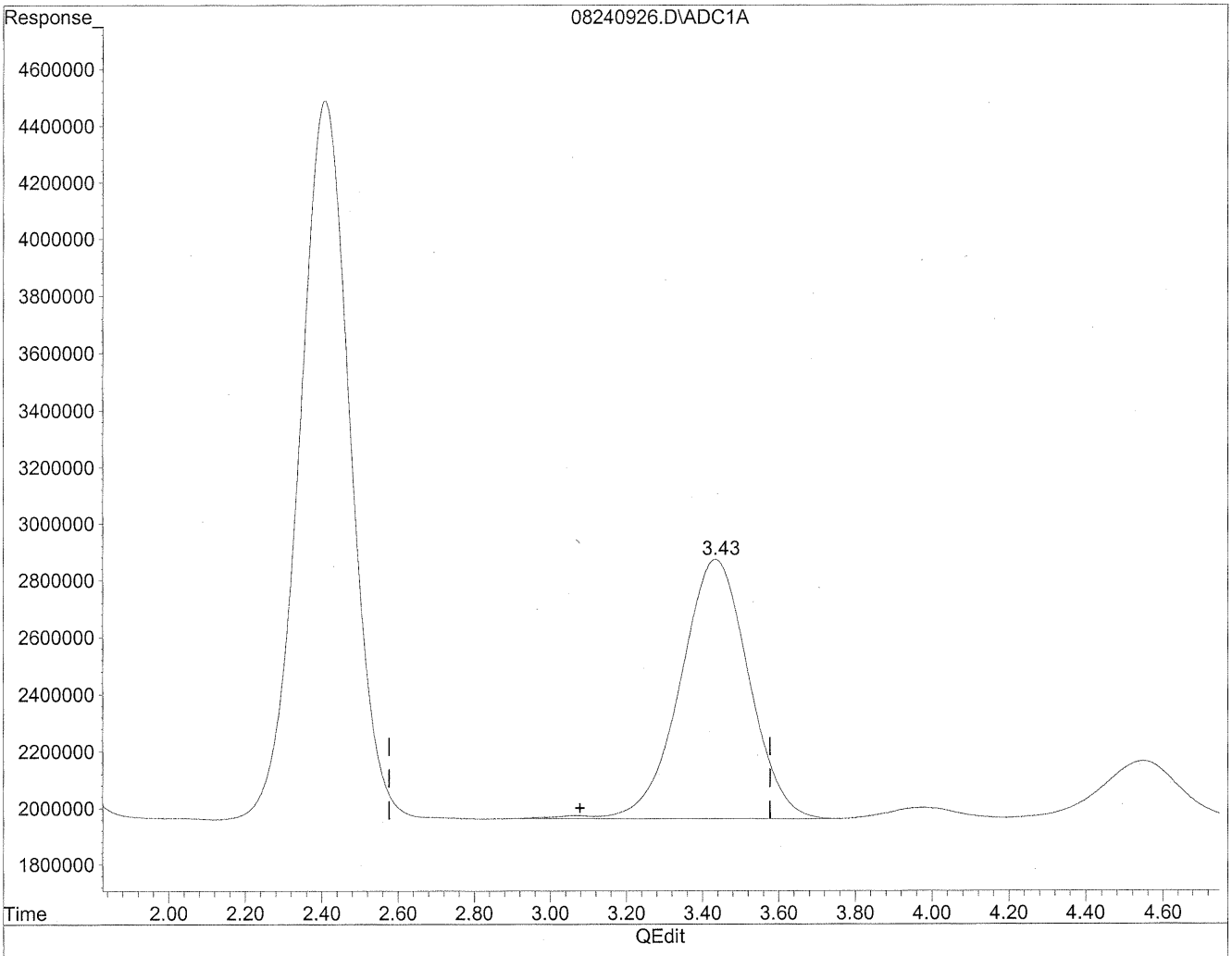
102

(+) = Expected Retention Time
08240926.D TO110709.M Sat Aug 29 13:37:35 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240926.D Vial: 24
Acq On : 24 Aug 2009 6:46 pm Operator: HC
Sample : P0902912-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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



(3) Propionaldehyde

3.43min 1065.886ng/ml

response 113724980

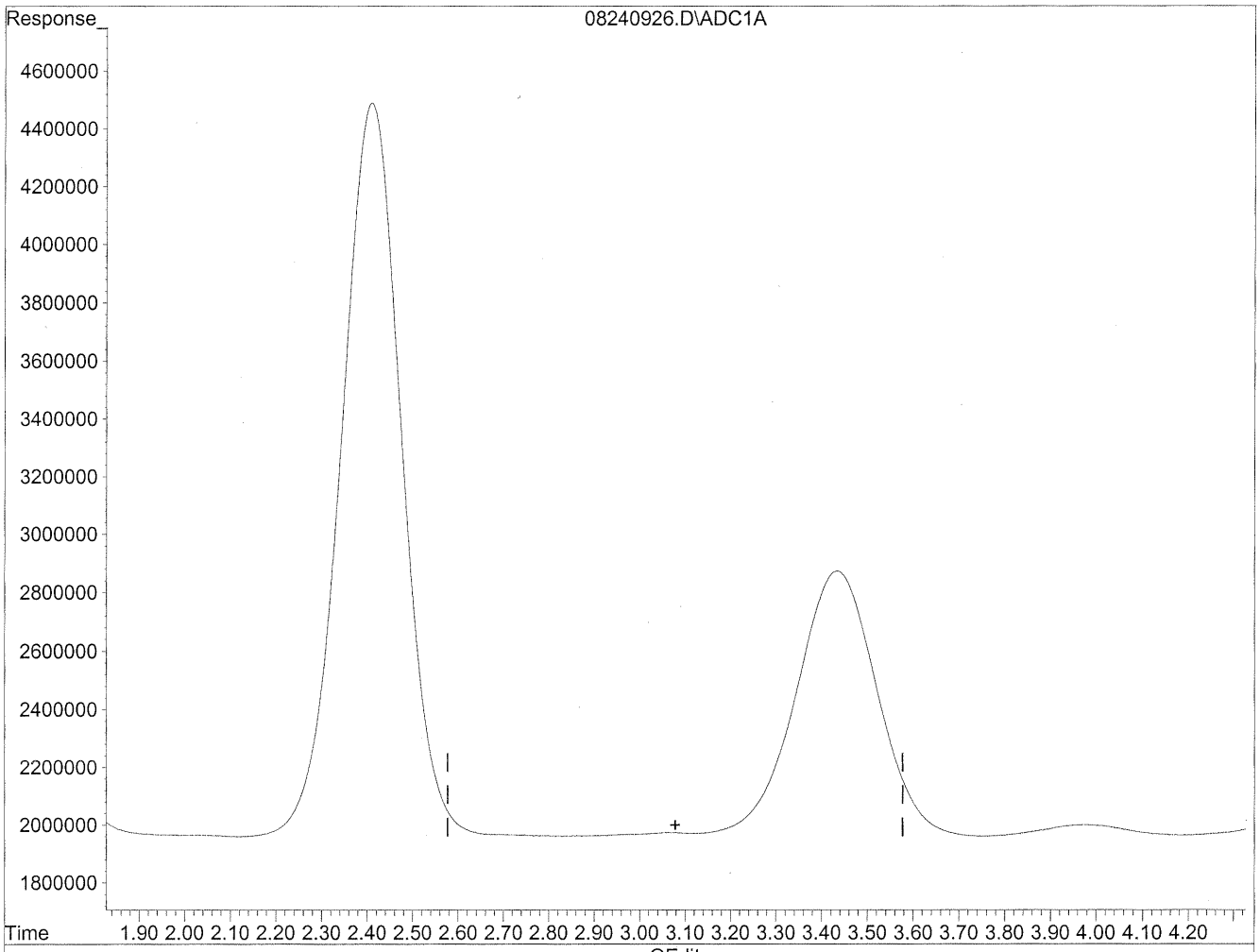
(+) = Expected Retention Time

08240926.D TO110709.M Sat Aug 29 13:37:38 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240926.D Vial: 24
Acq On : 24 Aug 2009 6:46 pm Operator: HC
Sample : P0902912-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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



QEdit

(3) Propionaldehyde

0.00min 0.000ng/ml d

response 0

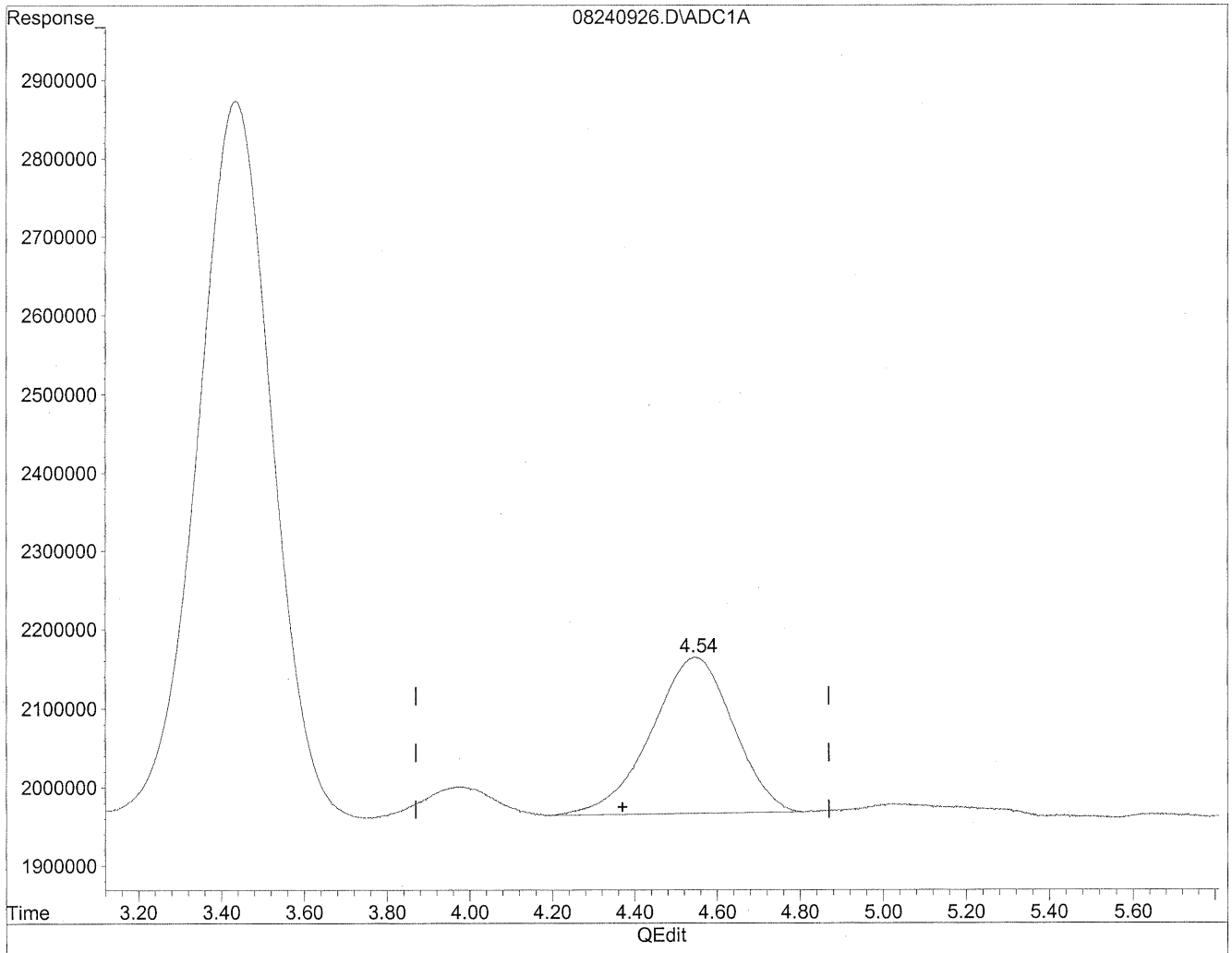
(+) = Expected Retention Time

08240926.D TO110709.M Sat Aug 29 13:37:40 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240926.D Vial: 24
Acq On : 24 Aug 2009 6:46 pm Operator: HC
Sample : P0902912-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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



(4) Crotonaldehyde

4.55min 280.462ng/ml

response 27321283

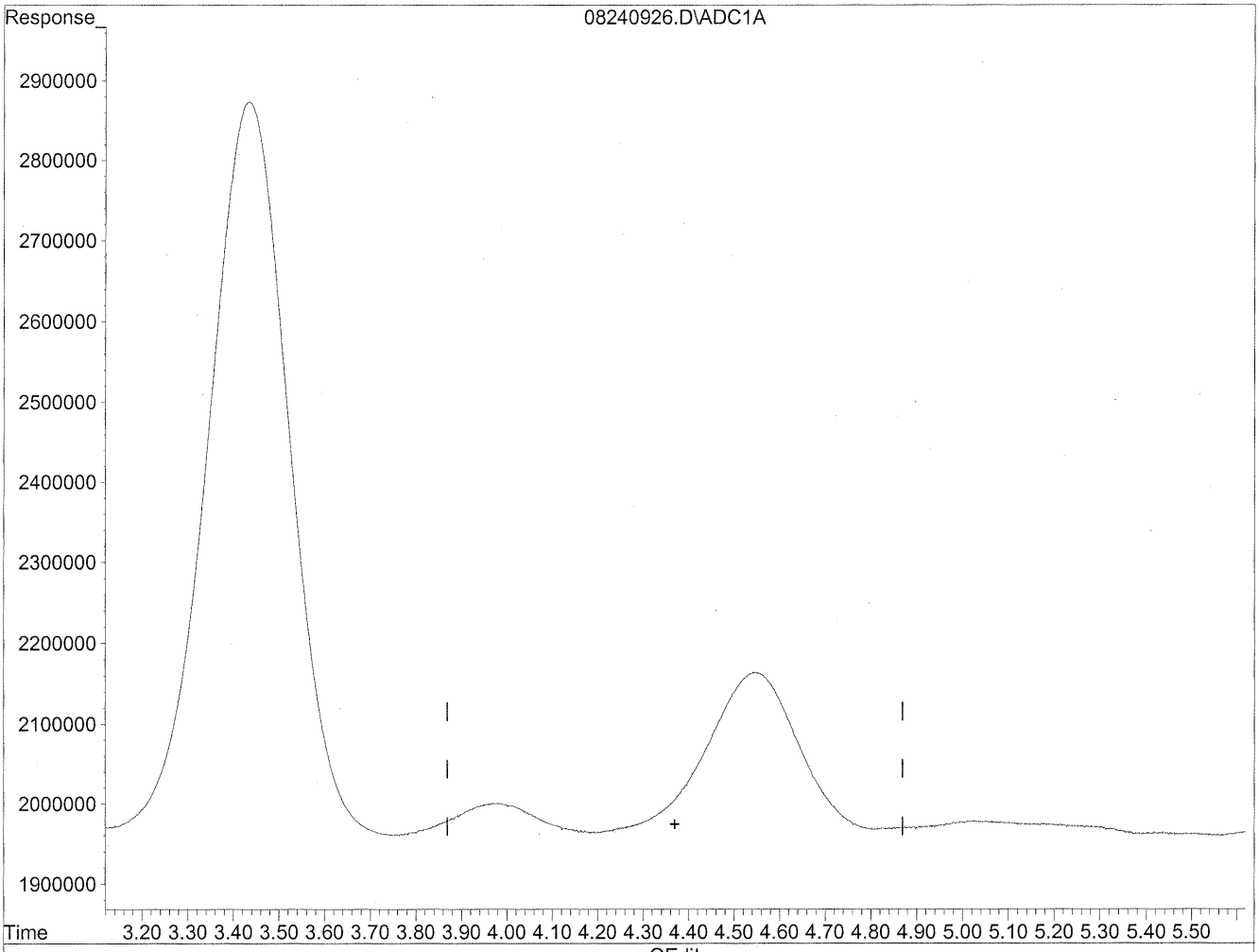
105

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240926.D Vial: 24
Acq On : 24 Aug 2009 6:46 pm Operator: HC
Sample : P0902912-004 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 19109 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



QEdit

(4) Crotonaldehyde

0.00min 0.000ng/ml d

response 0

106

(+) = Expected Retention Time

08240926.D TO110709.M Sat Aug 29 13:37:50 2009

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Sample ID: **102514**
Client Project ID: **16512**

CAS Project ID: P0902912
CAS Sample ID: P0902912-005

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: 8/20/09
Date Received: 8/21/09
Date Analyzed: 8/25/09
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	

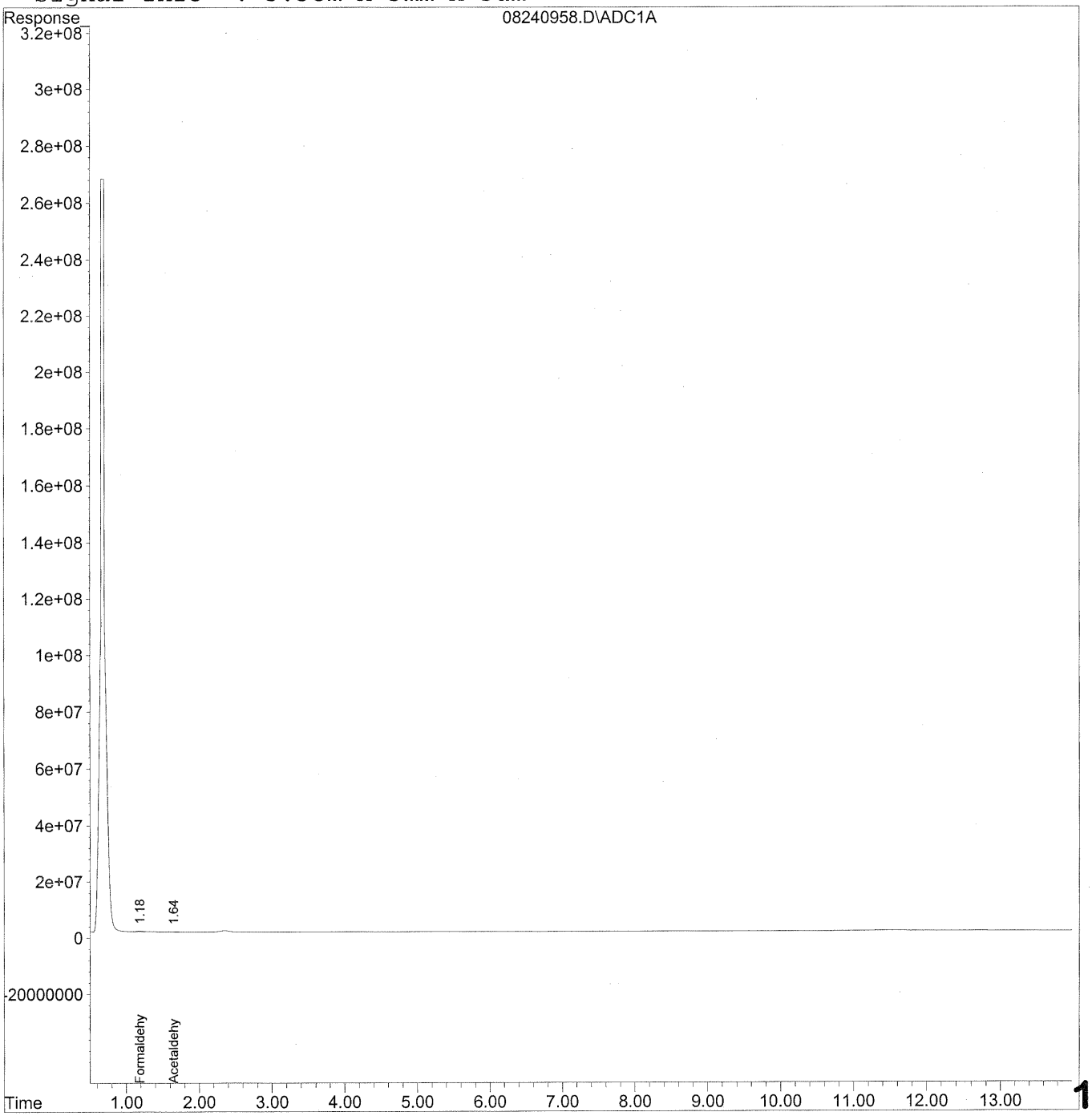
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
BC = Results reported are not blank corrected.
NA = Not applicable.

Verified By: Ro Date: 9/2/09

Data File : J:\LC01\DATA\TO11\2009_08\24\08240958.D Vial: 55
Acq On : 25 Aug 2009 2:47 am Operator: HC
Sample : P0902912-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 15:00 19109 Quant Results File: TO110709.RES

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



108

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\24\08240958.D Vial: 55
Acq On : 25 Aug 2009 2:47 am Operator: HC
Sample : P0902912-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 29 15:00 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

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

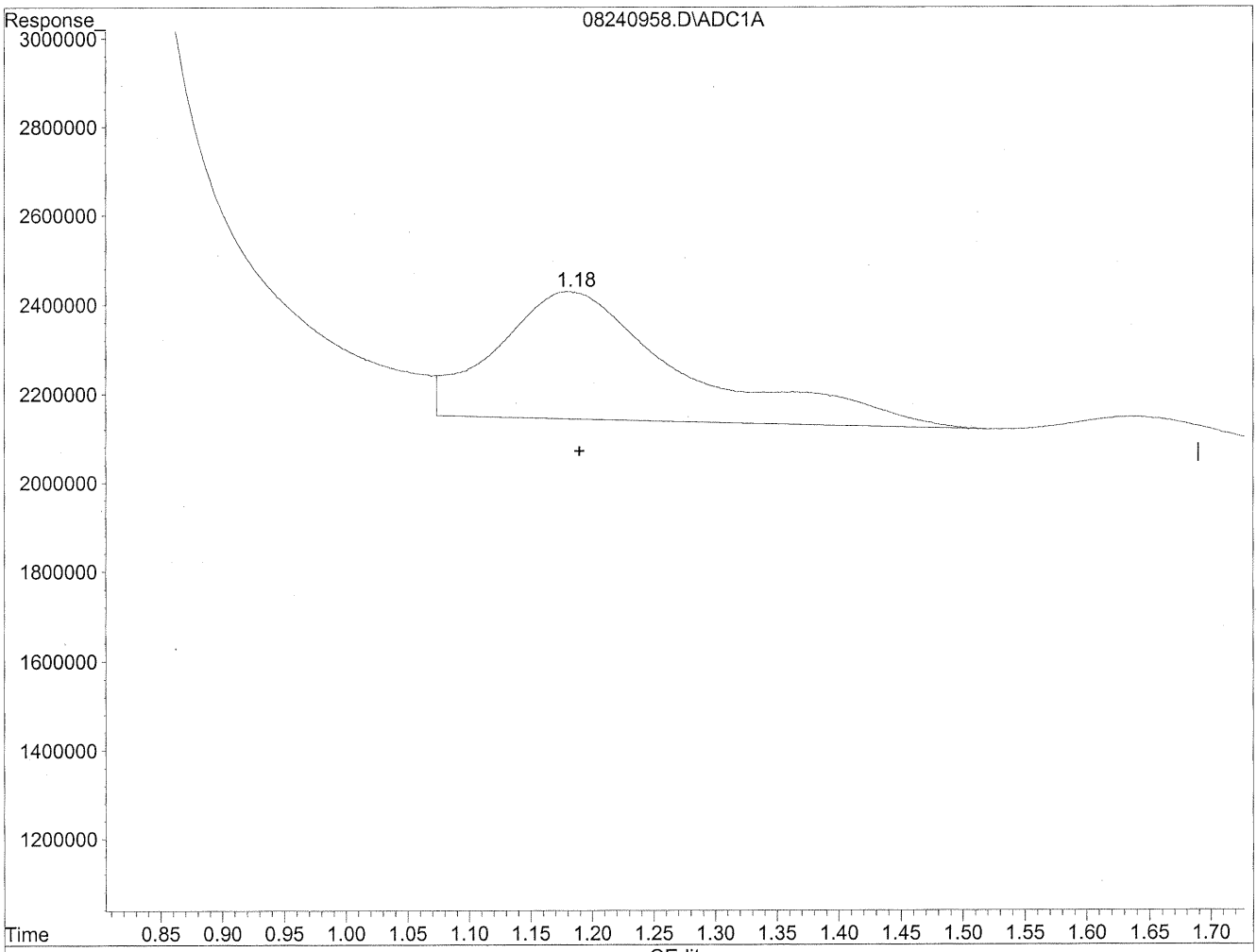
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.18	13242901	72.136	ng/mlm
2)	Acetaldehyde	1.64	3596091	25.645	ng/mlm
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	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/mld

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240958.D Vial: 55
Acq On : 25 Aug 2009 2:47 am Operator: HC
Sample : P0902912-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.18min 165.318ng/ml

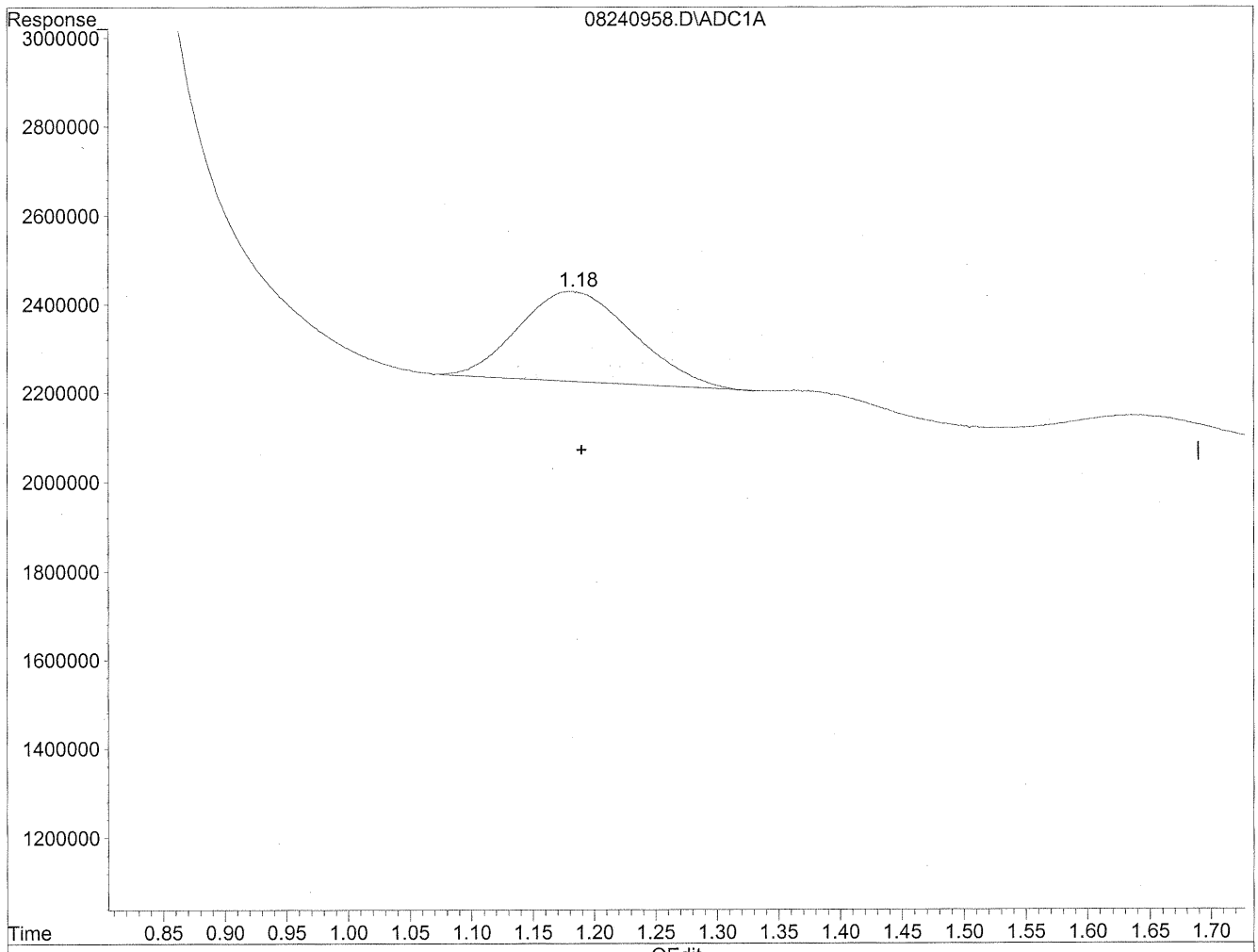
response 30349274

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240958.D Vial: 55
Acq On : 25 Aug 2009 2:47 am Operator: HC
Sample : P0902912-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.18min 72.136ng/ml m

response 13242901

HC 8/27/09

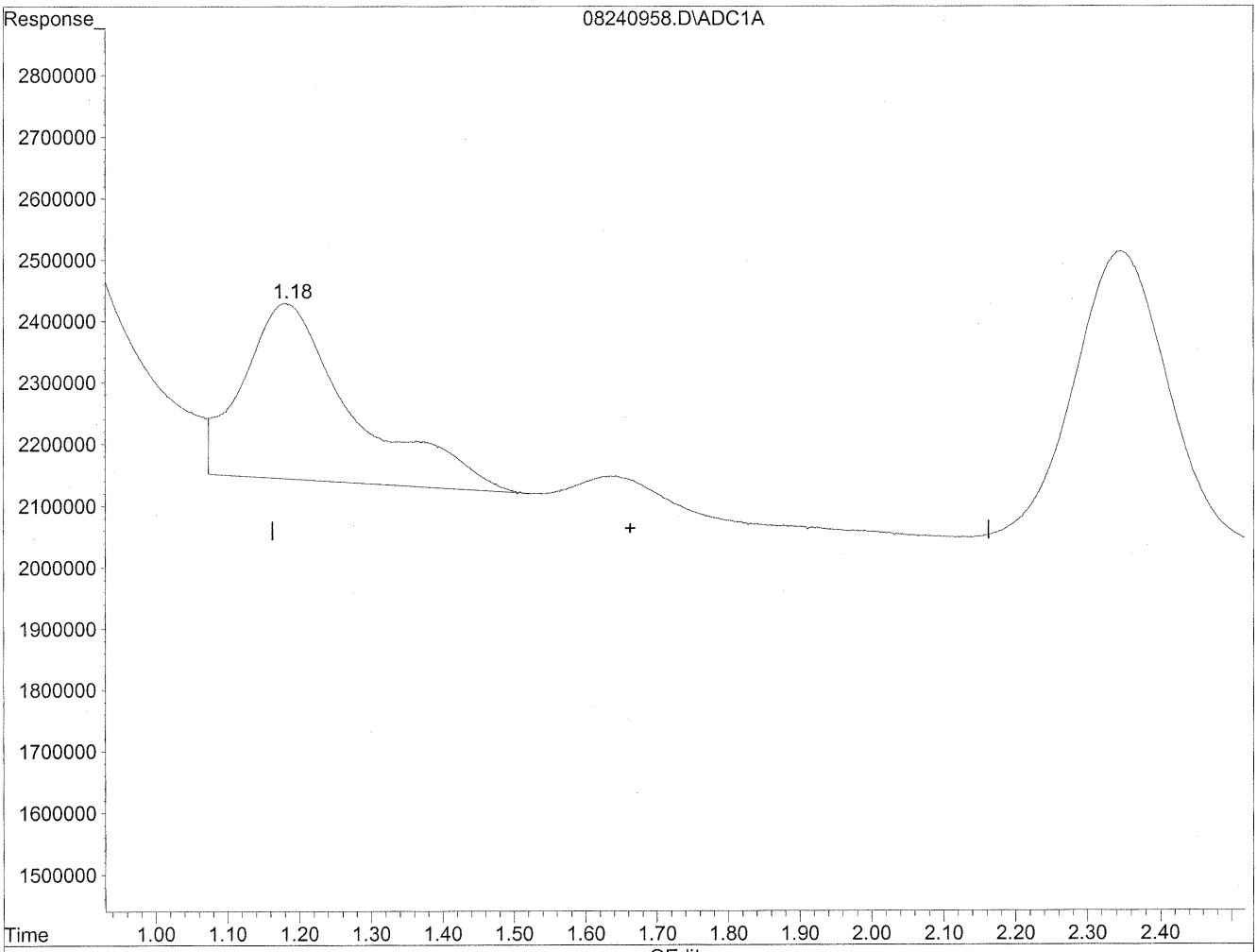
LC 8/27/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240958.D Vial: 55
Acq On : 25 Aug 2009 2:47 am Operator: HC
Sample : P0902912-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.18min 216.435ng/ml

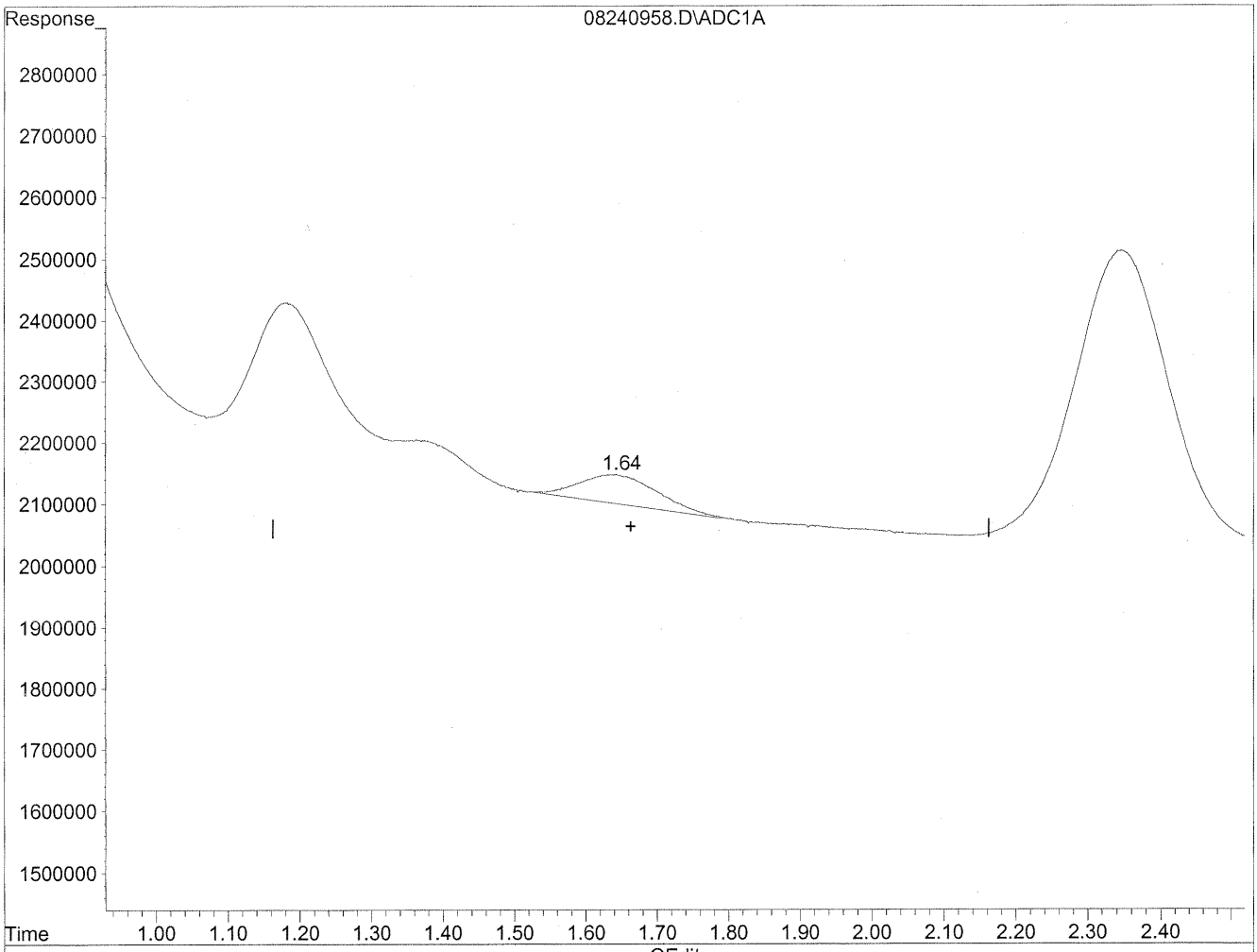
response 30349274

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240958.D Vial: 55
Acq On : 25 Aug 2009 2:47 am Operator: HC
Sample : P0902912-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.64min 25.645ng/ml m

response 3596091

HC 8/31/09

LW 9/3/09

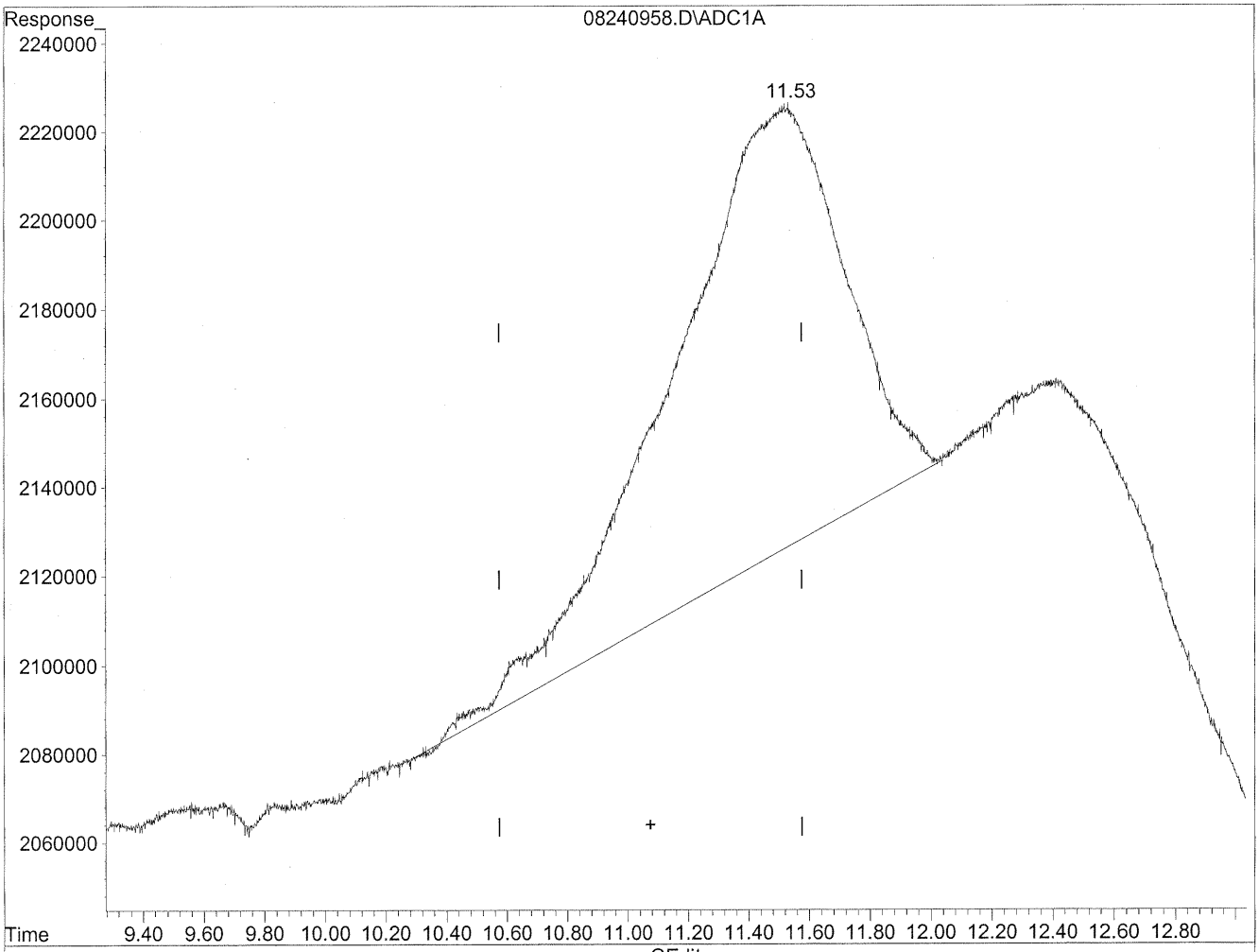
113

(+) = Expected Retention Time
08240958.D TO110709.M Sat Aug 29 15:00:47 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240958.D Vial: 55
Acq On : 25 Aug 2009 2:47 am Operator: HC
Sample : P0902912-005 front 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

11.52min 813.295ng/ml

response 39862329

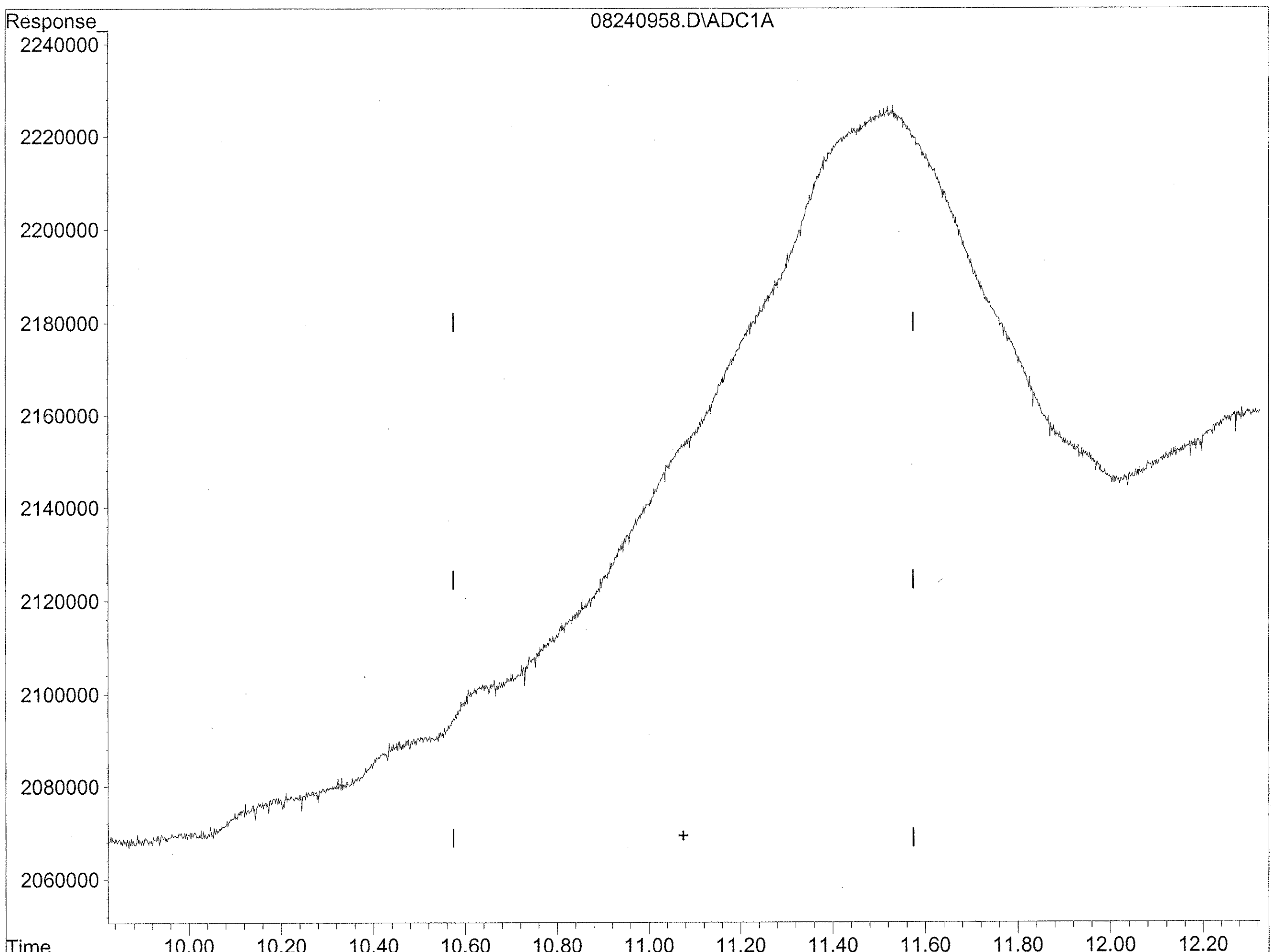
(+) = Expected Retention Time

08240958.D TO110709.M Sat Aug 29 15:00:51 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240958.D   Vial: 55
Acq On   : 25 Aug 2009   2:47 am   Operator: HC
Sample   : P0902912-005 front 1.0ml   Inst   : LC 01
Misc   :   Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Method   : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title   : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update   : Sat Aug 29 12:41:27 2009
Response via   : Multiple Level Calibration



QEdit

(12) 2,5-Dimethylbenzaldehyde

0.00min   0.000ng/ml d

response   0

HC 8/29/09 not real

LW 8/31/09

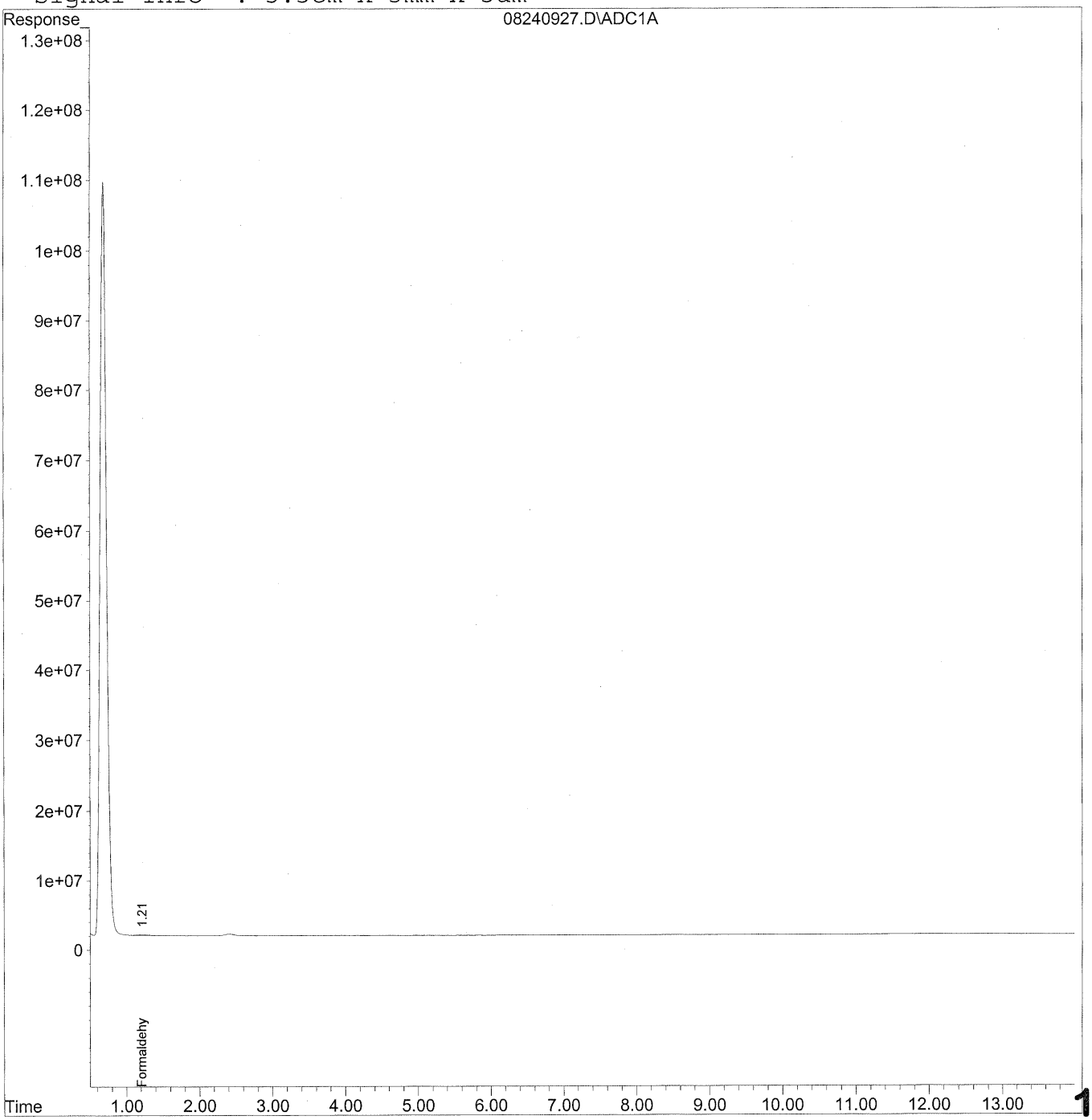
(+) = Expected Retention Time
08240958.D   TO110709.M   Sat Aug 29 15:00:54 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240927.D Vial: 25
Acq On : 24 Aug 2009 7:01 pm Operator: HC
Sample : P0902912-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 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 : Tue Aug 25 10:21:57 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\24\08240927.D Vial: 25
Acq On : 24 Aug 2009 7:01 pm Operator: HC
Sample : P0902912-005 back 1.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 14:51 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 : Tue Aug 25 10:21:57 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	1.21	12887504	70.200	ng/ml
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

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

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

CAS Project ID: P0902912
CAS Sample ID: P090824-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: 08/24/09
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.

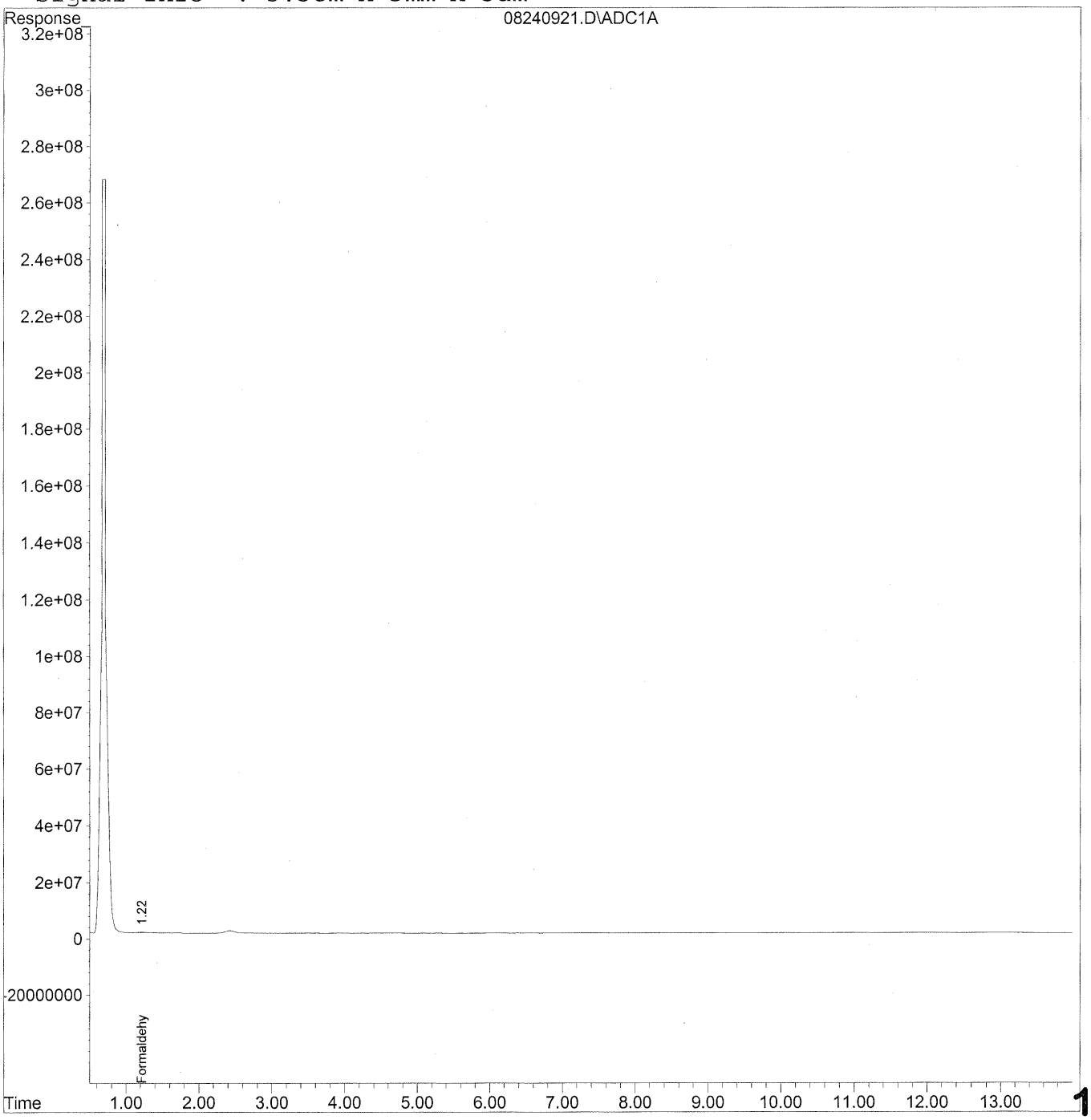
Verified By: RC Date: 9/2/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240921.D Vial: 19
Acq On : 24 Aug 2009 5:31 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 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 : 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\24\08240921.D Vial: 19
Acq On : 24 Aug 2009 5:31 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 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 : 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

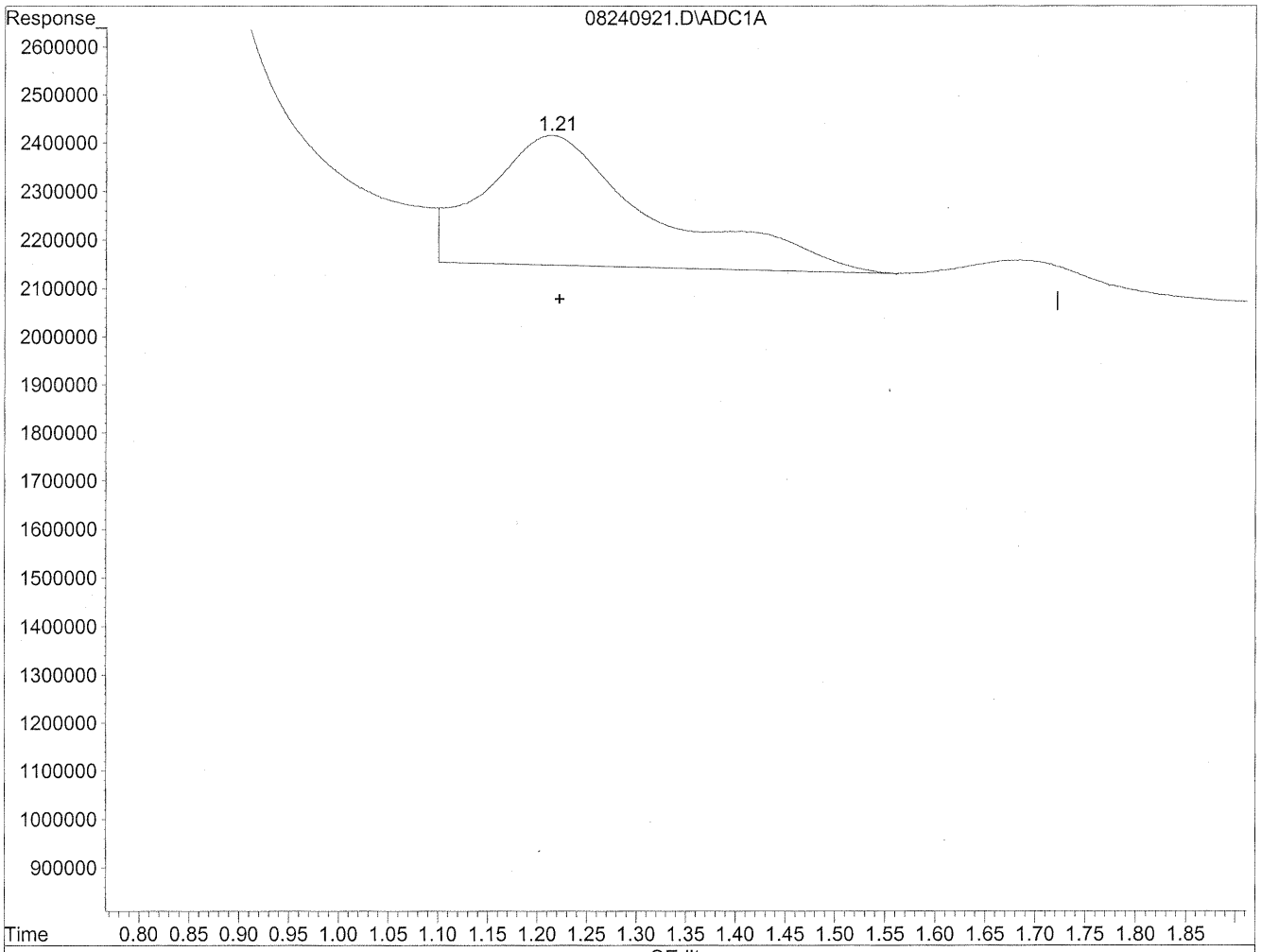
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	11322887	61.678	ng/mlm
2)	Acetaldehyde	0.00	0	N.D.	ng/ml
3)	Propionaldehyde	0.00	0	N.D.	ng/ml
4)	Crotonaldehyde	0.00	0	N.D.	ng/ml
5)	Butyraldehyde	0.00	0	N.D.	ng/ml
6)	Benzaldehyde	0.00	0	N.D.	ng/ml
7)	Isovaleraldehyde	0.00	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

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240921.D Vial: 19
Acq On : 24 Aug 2009 5:31 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 13:14 19109 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



QEdit

(1) Formaldehyde

1.21min 171.768ng/ml

response 31533394

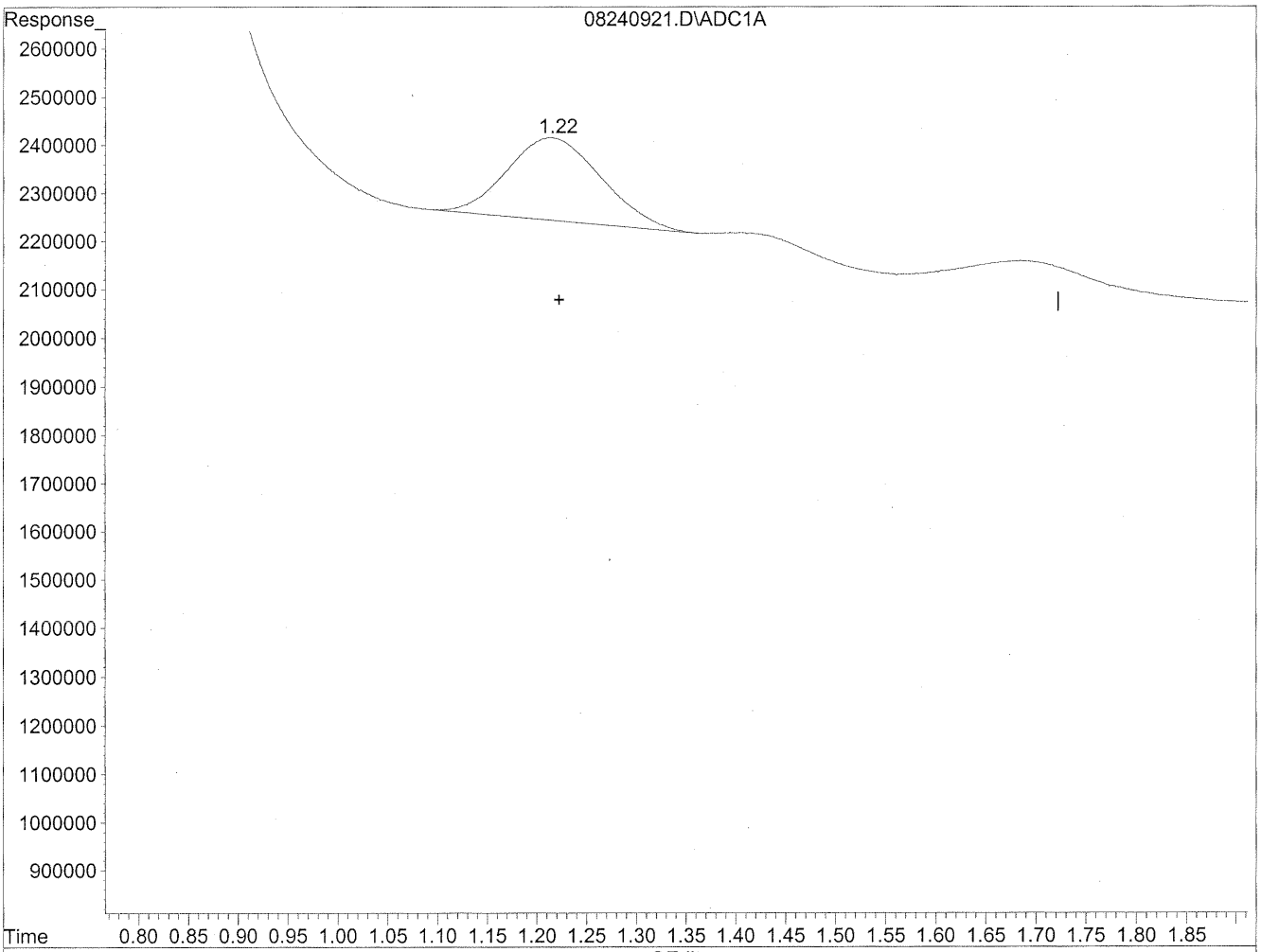
121

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240921.D Vial: 19
Acq On : 24 Aug 2009 5:31 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 13:14 19109 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



QEdit

(1) Formaldehyde

1.22min 61.678ng/ml m

response 11322887

8/29/09

122

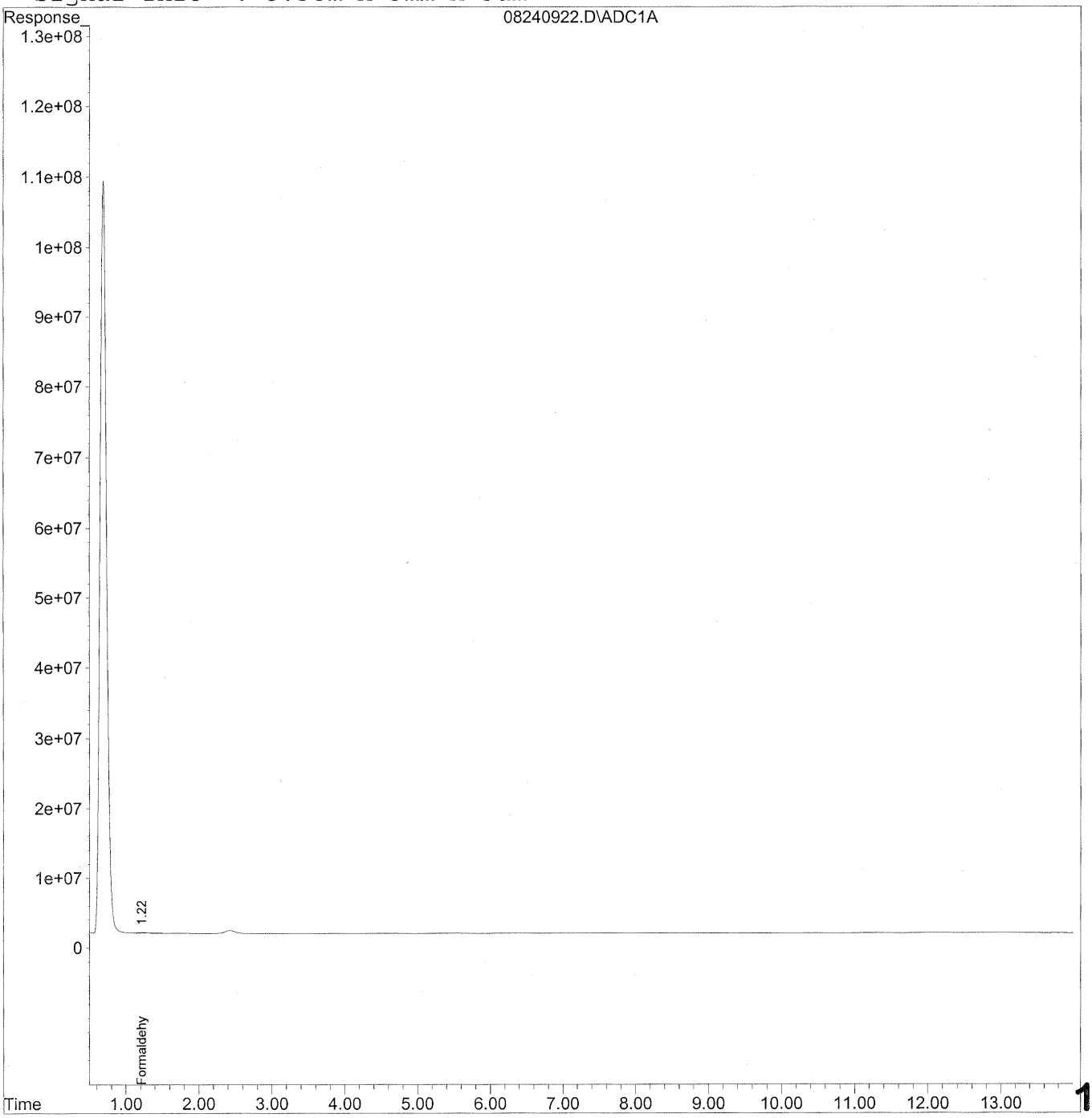
(+) = Expected Retention Time
08240921.D TO110709.M Sat Aug 29 13:15:02 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240922.D Vial: 20
Acq On : 24 Aug 2009 5:46 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 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 : 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



123

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\24\08240922.D Vial: 20
Acq On : 24 Aug 2009 5:46 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 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 : 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	1.21	14557359	79.296	ng/ml
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

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

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

CAS Project ID: P0902912
CAS Sample ID: P090825-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: 08/25/09
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.

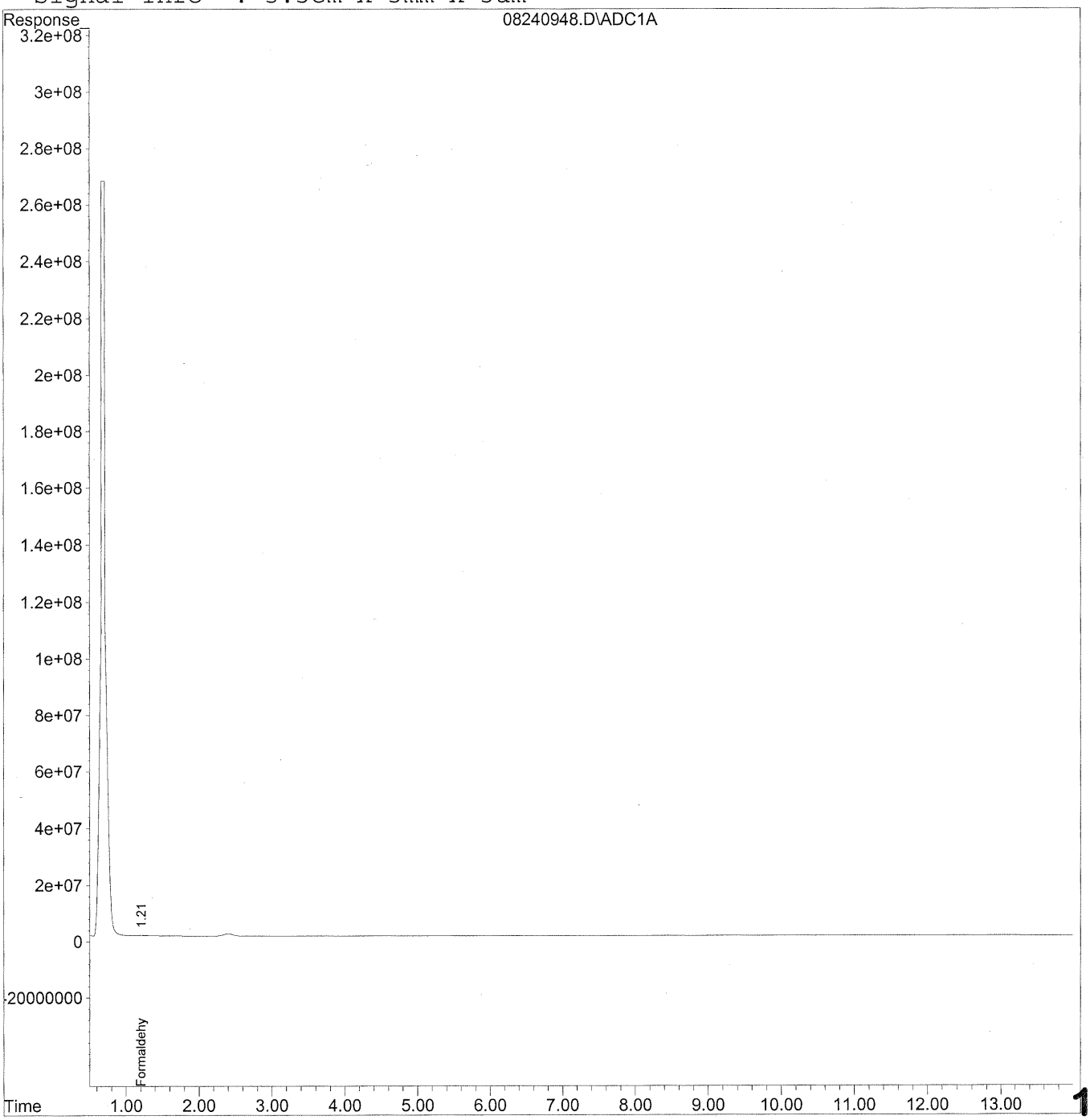
Verified By: Rc Date: 9/2/09

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240948.D Vial: 45
Acq On : 25 Aug 2009 12:17 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 13:17 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : 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



126

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\24\08240948.D Vial: 45
Acq On : 25 Aug 2009 12:17 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 13:17 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

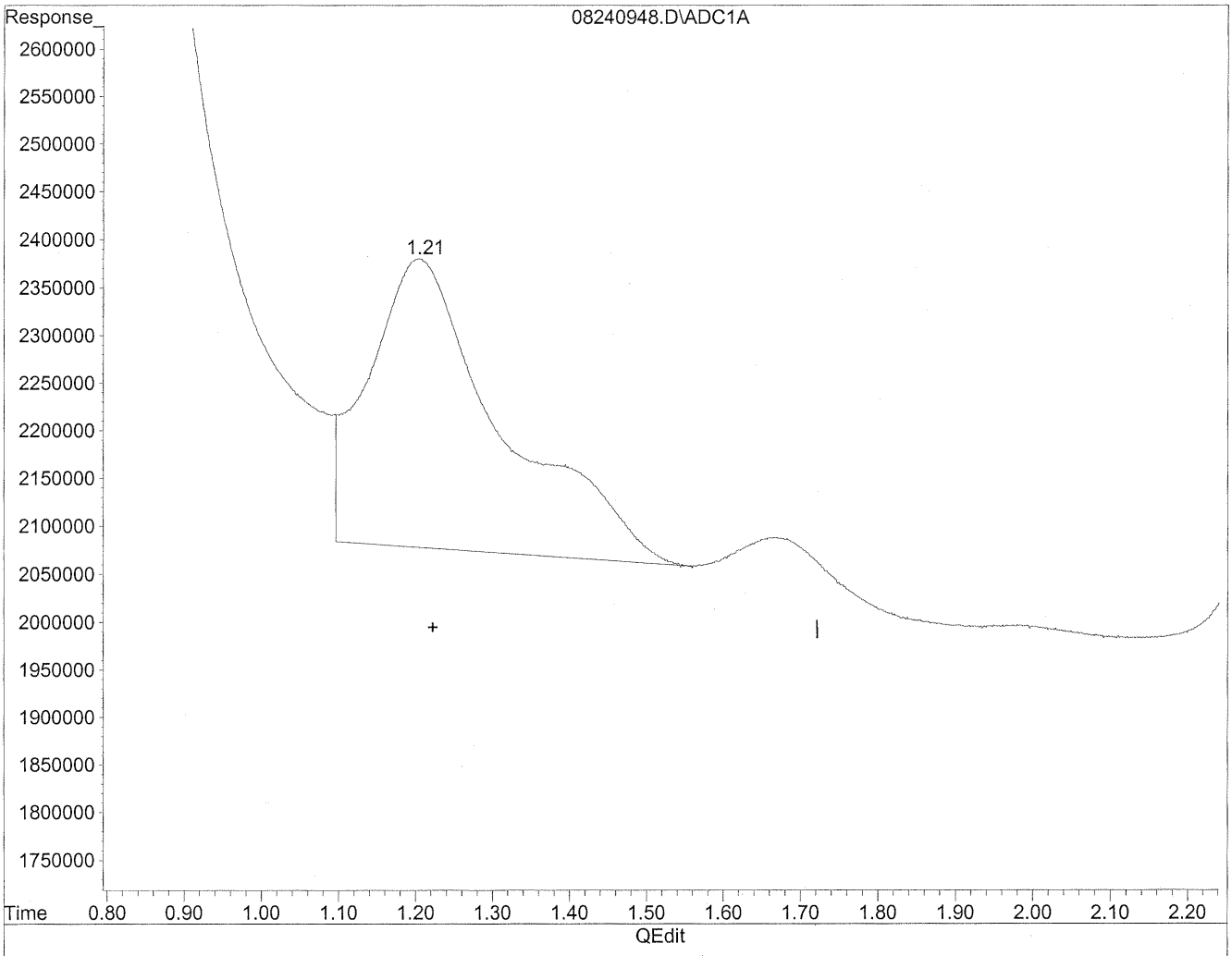
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.21	12298739	66.993	ng/mlm
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\24\08240948.D Vial: 45
Acq On : 25 Aug 2009 12:17 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 13:16 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.20min 197.992ng/ml

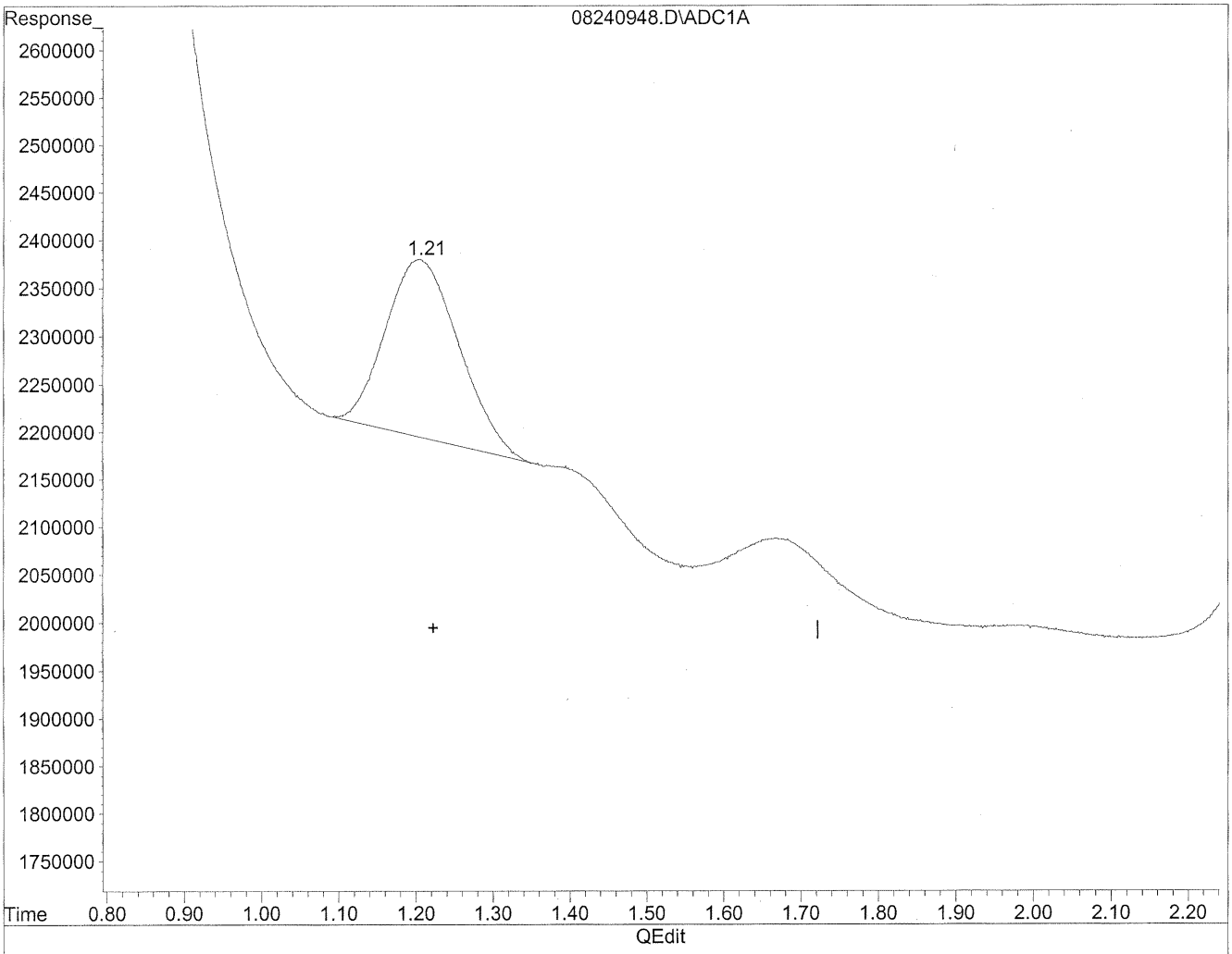
response 36347573

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240948.D Vial: 45
Acq On : 25 Aug 2009 12:17 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 13:16 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Sat Aug 29 12:41:27 2009
Response via : Multiple Level Calibration



(1) Formaldehyde

1.21min 66.993ng/ml m

response 12298739

pc 8/29/09 LC

LB 8/31/09

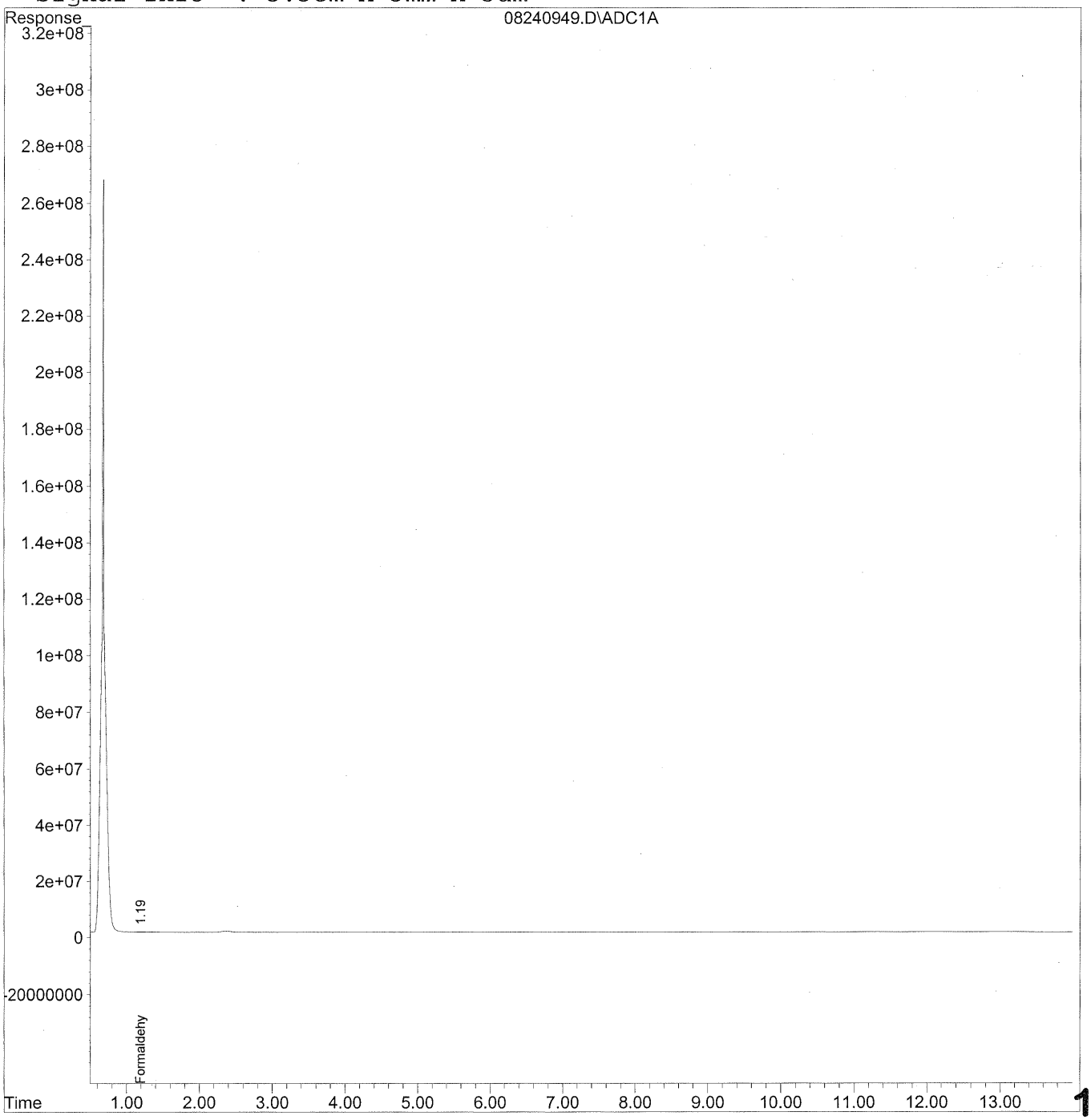
129

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2009_08\24\08240949.D Vial: 46
Acq On : 25 Aug 2009 12:32 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 13:17 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : 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\24\08240949.D Vial: 46
Acq On : 25 Aug 2009 12:32 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 13:17 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : 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	1.19	16684741	90.885	ng/ml
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

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%	4293221	0.43%	3079204	0.70%
50ng/ml TO11A Std	9305088	4.78%	7389770	7.24%	5442713	6.66%	5754474	6.32%	4119144	3.64%	2732056	10.66%
100ng/ml TO11A St	18283557	0.51%	13784712	1.44%	10870707	0.86%	9346475	1.91%	8839595	0.81%	7282249	5.41%
100ng/ml TO11A St	18449443	0.39%	14434553	3.21%	11389784	3.88%	9814490	3.00%	9432197	5.84%	6706722	2.92%
100ng/ml TO11A St	18400032	0.12%	13737532	1.77%	10633406	3.02%	9424529	1.09%	8463028	5.03%	6735919	2.50%
500ng/ml TO11A St	91593554	0.39%	70468869	0.90%	53468174	1.20%	47866960	1.26%	43271557	0.62%	32616313	2.91%
500ng/ml TO11A St	90711575	0.57%	69140255	1.00%	52850412	0.03%	47584179	0.66%	43677338	0.31%	34085310	1.46%
500ng/ml TO11A St	91399555	0.18%	69908753	0.10%	52190620	1.22%	46362546	1.92%	43673214	0.30%	34084716	1.46%
1500ng/ml TO11A S	275380897	0.26%	209374751	0.16%	159030091	0.21%	143227783	1.11%	134132687	1.08%	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%	706170560	0.05%	539067854	0.39%	476268543	0.19%	446392739	0.21%	328286106	0.04%
5000ng/ml TO11A S	925768000	0.17%	708552415	0.38%	540133923	0.59%	477844499	0.52%	446568052	0.25%	328413551	0.08%
5000ng/ml TO11A S	918424042	0.62%	702791887	0.43%	531675082	0.98%	471954575	0.72%	443441833	0.45%	327762901	0.12%
10000ng/ml TO11A	1908653125	0.62%	1450154617	0.67%	1099941045	0.36%	972691462	0.37%	910896701	0.36%	668462127	0.28%
10000ng/ml TO11A	1905913073	0.48%	1446499891	0.41%	1098837646	0.26%	971357788	0.23%	911328243	0.41%	669128969	0.38%
10000ng/ml TO11A	1875917434	1.10%	1425028469	1.08%	1089338811	0.61%	963283335	0.60%	900561239	0.78%	662238443	0.66%

AVERAGE RESPONSE FACTOR

HC 7/29/09

Method:
Analyst:

CALIBRATION F

Calibration Level	Isovaler-Aldehyde	% rpd	Valer-Aldehyde	% rpd	o-Tolu-Aldehyde	% rpd	m,p-Tolu-Aldehyde	% rpd	Hex-Aldehyde	% rpd	2,5-Dimethl benz-Aldehyde	% rpd
50ng/ml TO11A Std	4167653	7.13%	3532734	7.15%	338/183	22.94%	5445142	7.87%	3244418	5.31%	2546144	7.64%
50ng/ml TO11A Std	4002738	2.89%	4025564	5.81%	2461625	10.65%	489/087	2.98%	3295067	3.83%	2605446	5.49%
50ng/ml TO11A Std	3500271	10.02%	3855749	1.34%	2416389	12.29%	4801019	4.89%	3739368	9.14%	3118537	13.13%
100ng/ml TO11A St	7487274	5.83%	7060988	8.24%	5548699	2.73%	109/9457	1.36%	6702769	5.76%	5399082	9.13%
100ng/ml TO11A St	8338385	4.88%	811/341	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%	111/7259	0.42%	6920120	2.70%	4707951	4.84%
500ng/ml TO11A St	37944016	3.60%	3557/509	1.84%	29317615	1.49%	5327/975	1.62%	32888440	1.80%	23823948	2.62%
500ng/ml TO11A St	40968120	4.08%	36648075	1.12%	29793454	0.11%	54514161	0.67%	31855201	1.40%	22510750	3.03%
500ng/ml TO11A St	39175205	0.48%	36501988	0.72%	30169058	1.37%	54668231	0.95%	32179520	0.40%	23309464	0.41%
1500ng/ml TO11A S	115866442	0.09%	107104204	0.36%	86339652	0.42%	162946532	1.14%	98895406	0.29%	69932636	0.37%
1500ng/ml TO11A S	116723586	0.83%	107107592	0.37%	85940120	0.88%	161094009	0.01%	98090122	0.53%	6887/3541	1.15%
1500ng/ml TO11A S	114690000	0.92%	105937177	0.73%	87824227	1.30%	159292531	1.13%	98846718	0.24%	70224395	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%	35967/6615	0.47%	30007/384	0.48%	547211501	0.27%	333701808	0.31%	237108293	0.30%
5000ng/ml TO11A S	386992833	0.28%	356464469	0.43%	29737/4461	0.43%	544331756	0.26%	332038452	0.19%	23642820/	0.01%
10000ng/ml TO11A	790328317	0.44%	730218673	0.36%	60820827/6	0.16%	1111180147	0.26%	67351680/	0.25%	47846094/	0.27%
10000ng/ml TO11A	788026190	0.15%	729839210	0.31%	6103262/38	0.50%	11132098/0	0.45%	681915785	0.99%	484763918	1.04%
10000ng/ml TO11A	782256804	0.59%	722749626	0.67%	6032565/9	0.66%	11003845/3	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							

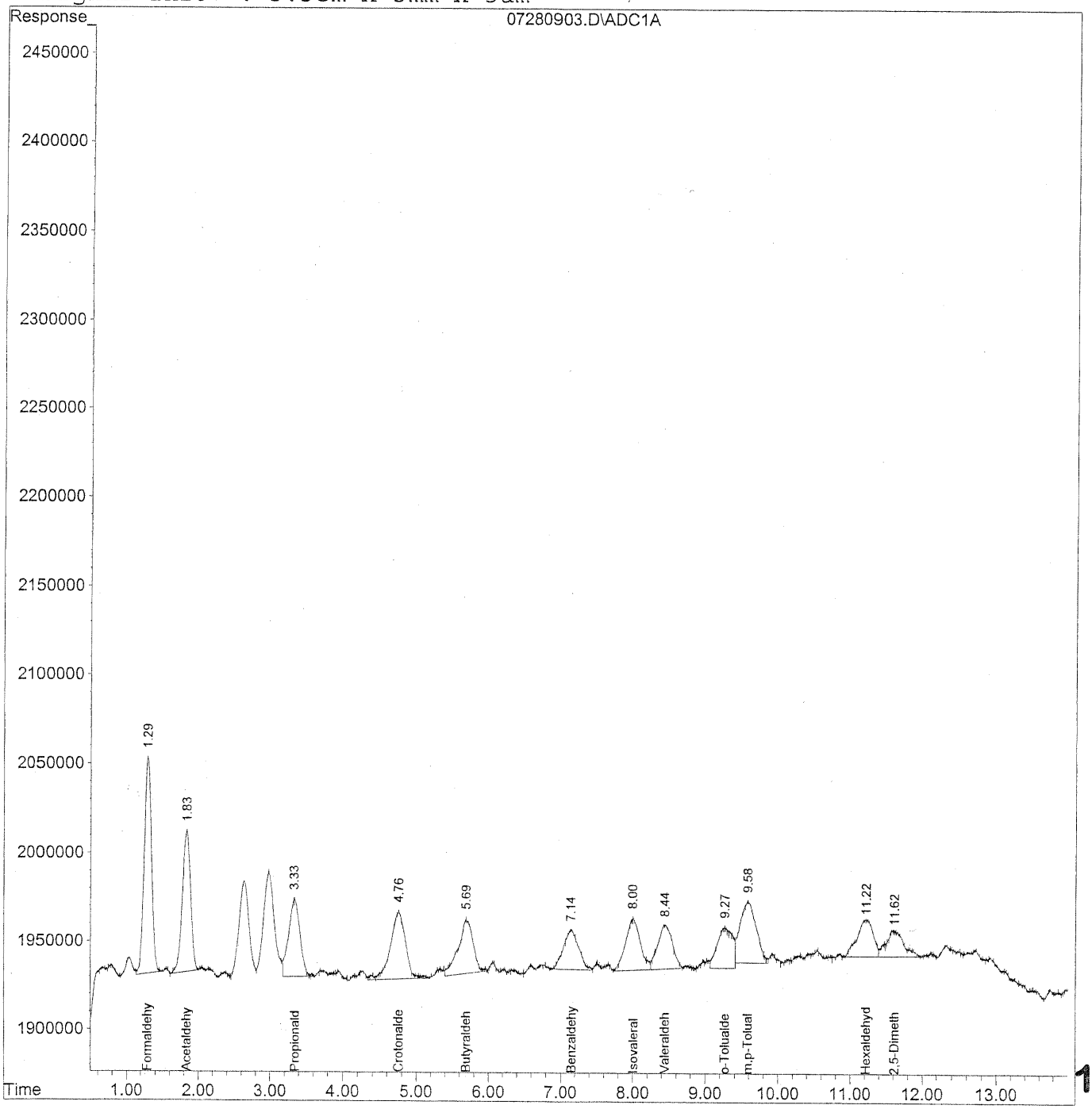
HC
7/29/04

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280903.D Vial: 3
Acq On : 28 Jul 2009 9:09 am Operator: HC
Sample : 50ng/ml TO11A Std S21-07270908 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:22 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

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



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

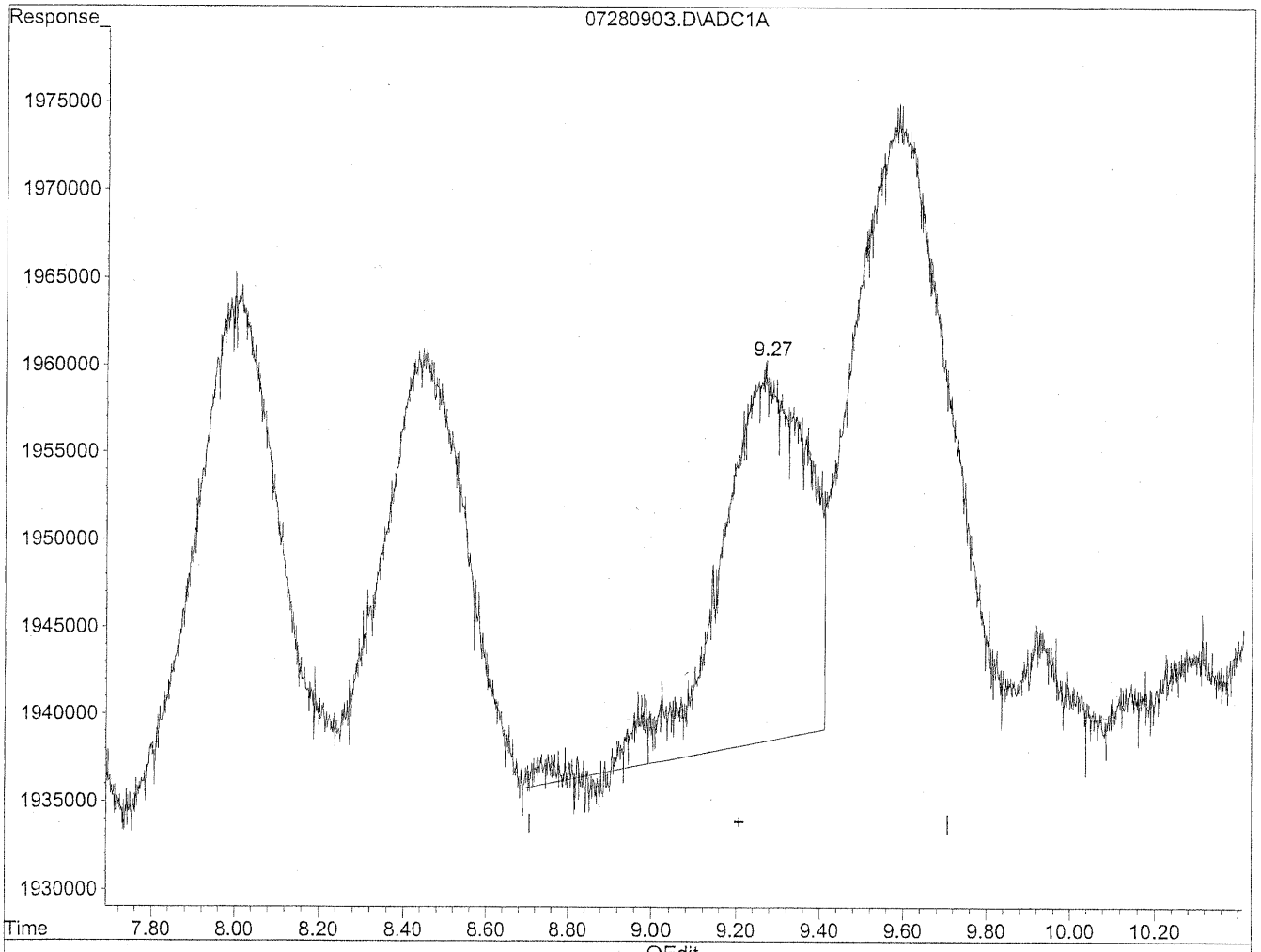
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.29	8477013	48.277	ng/ml
2)	Acetaldehyde	1.83	6307171	46.755	ng/ml
3)	Propionaldehyde	3.34	4892636	47.596	ng/ml
4)	Crotonaldehyde	4.76	5507079	49.813	ng/ml
5)	Butyraldehyde	5.70	4412295	54.828	ng/ml
6)	Benzaldehyde	7.15	3362429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

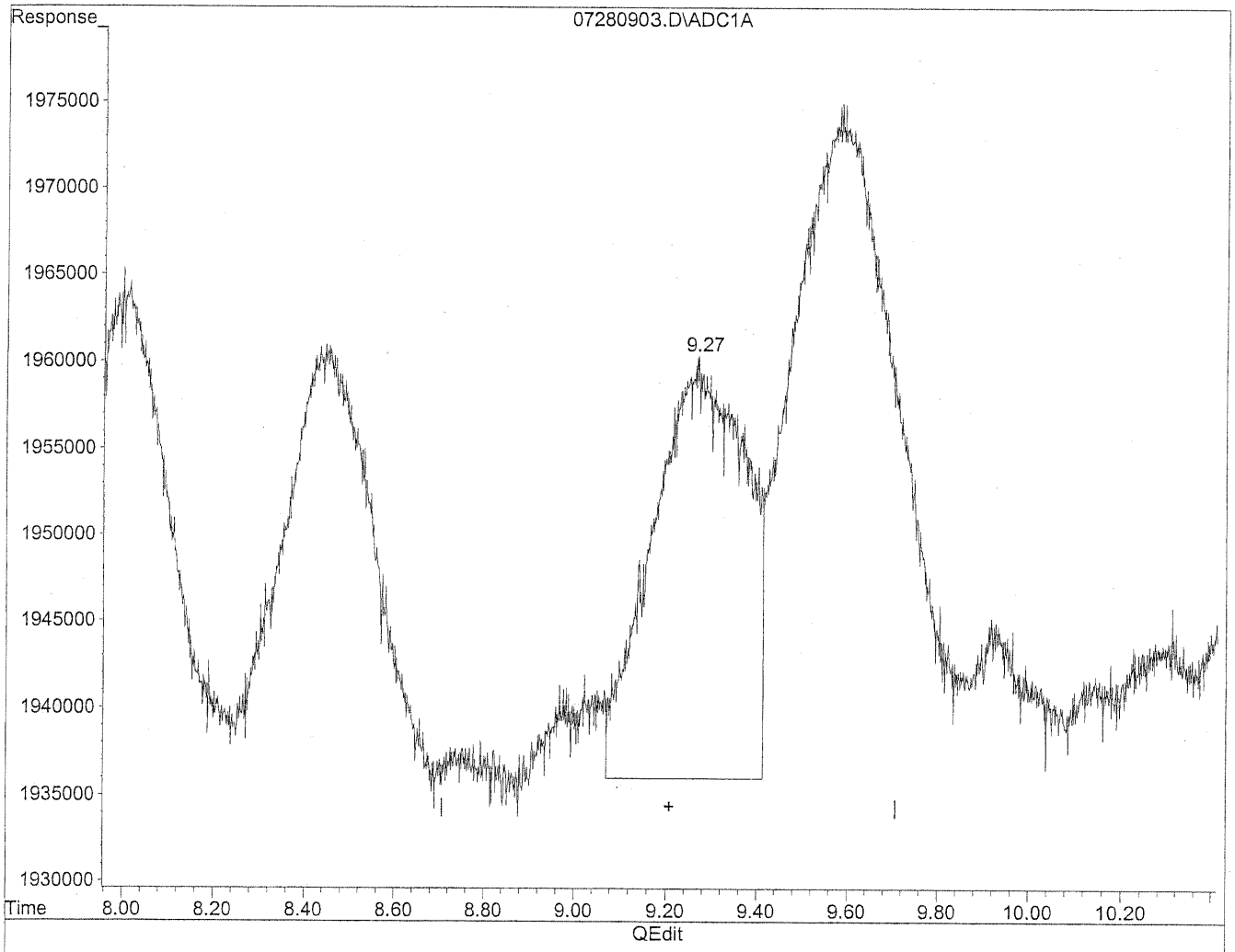
response 3109441

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280903.D Vial: 3
Acq On : 28 Jul 2009 9:09 am Operator: HC
Sample : 50ng/ml TO11A Std S21-07270908 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:22 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 15:29:52 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.27min 62.877ng/ml m

response 3387183

HL
Injector
1C

KF 7/29/09

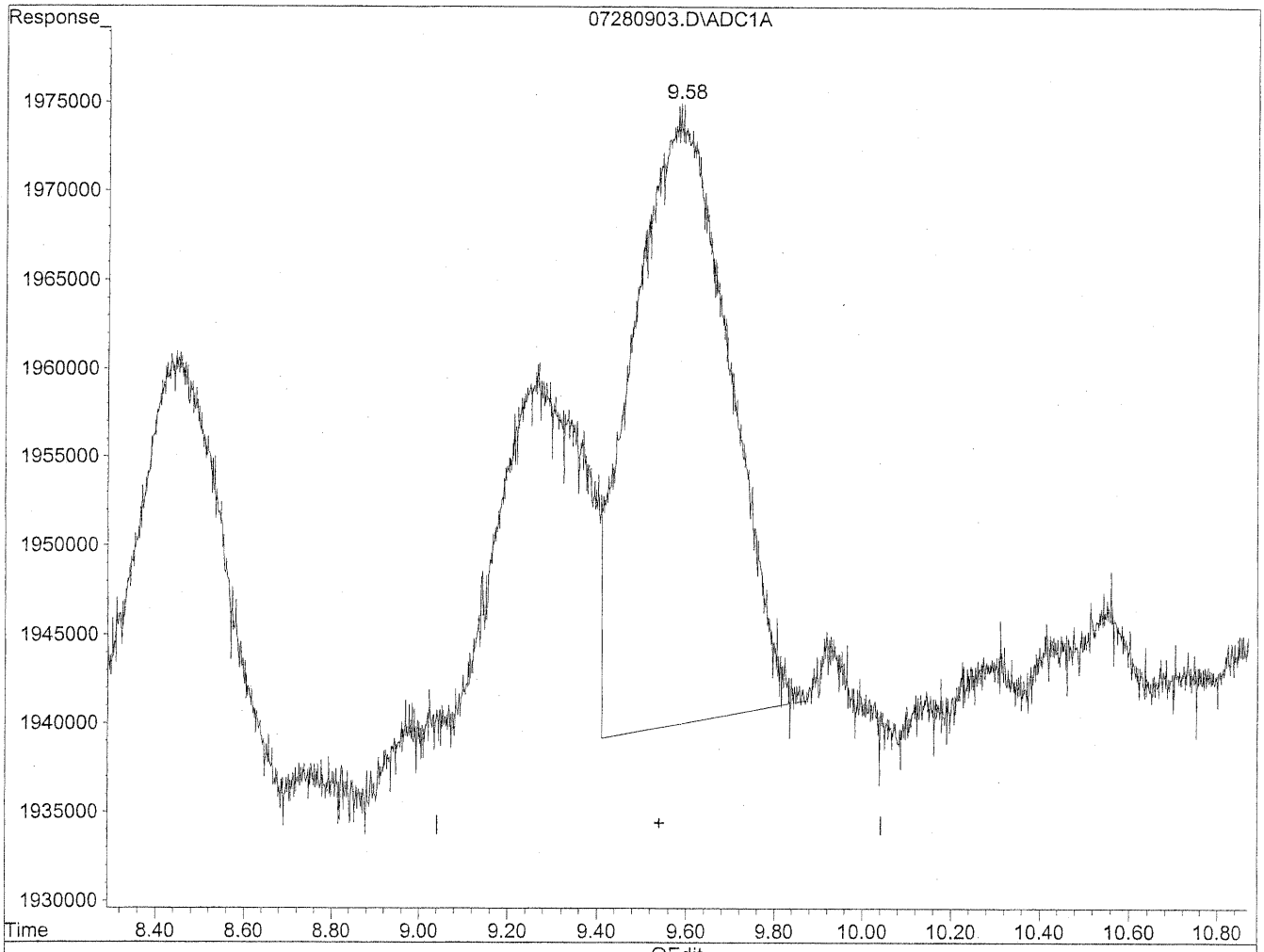
144

(+) = 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

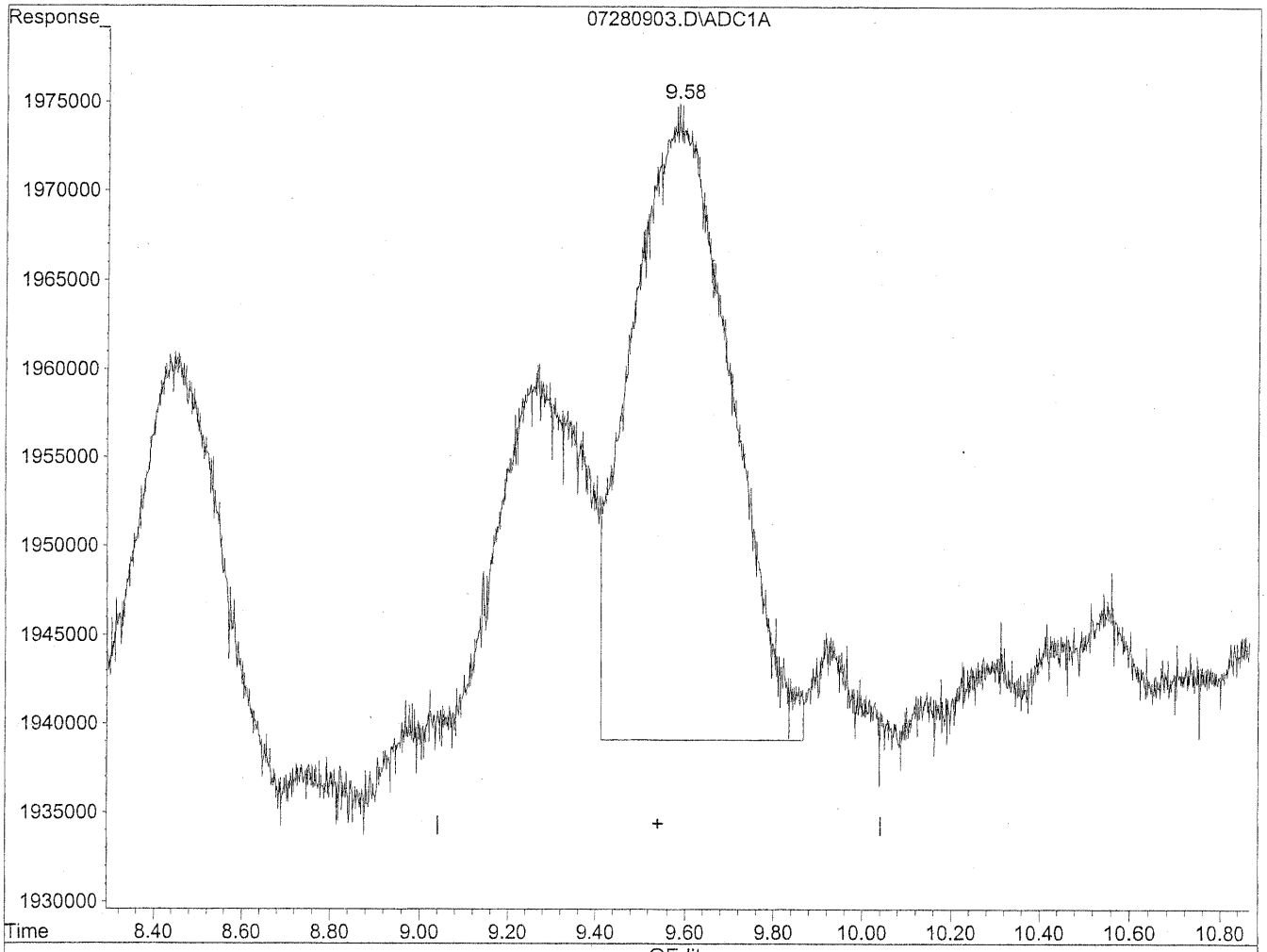
response 5147699

(+) = 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

KR 7/29/09

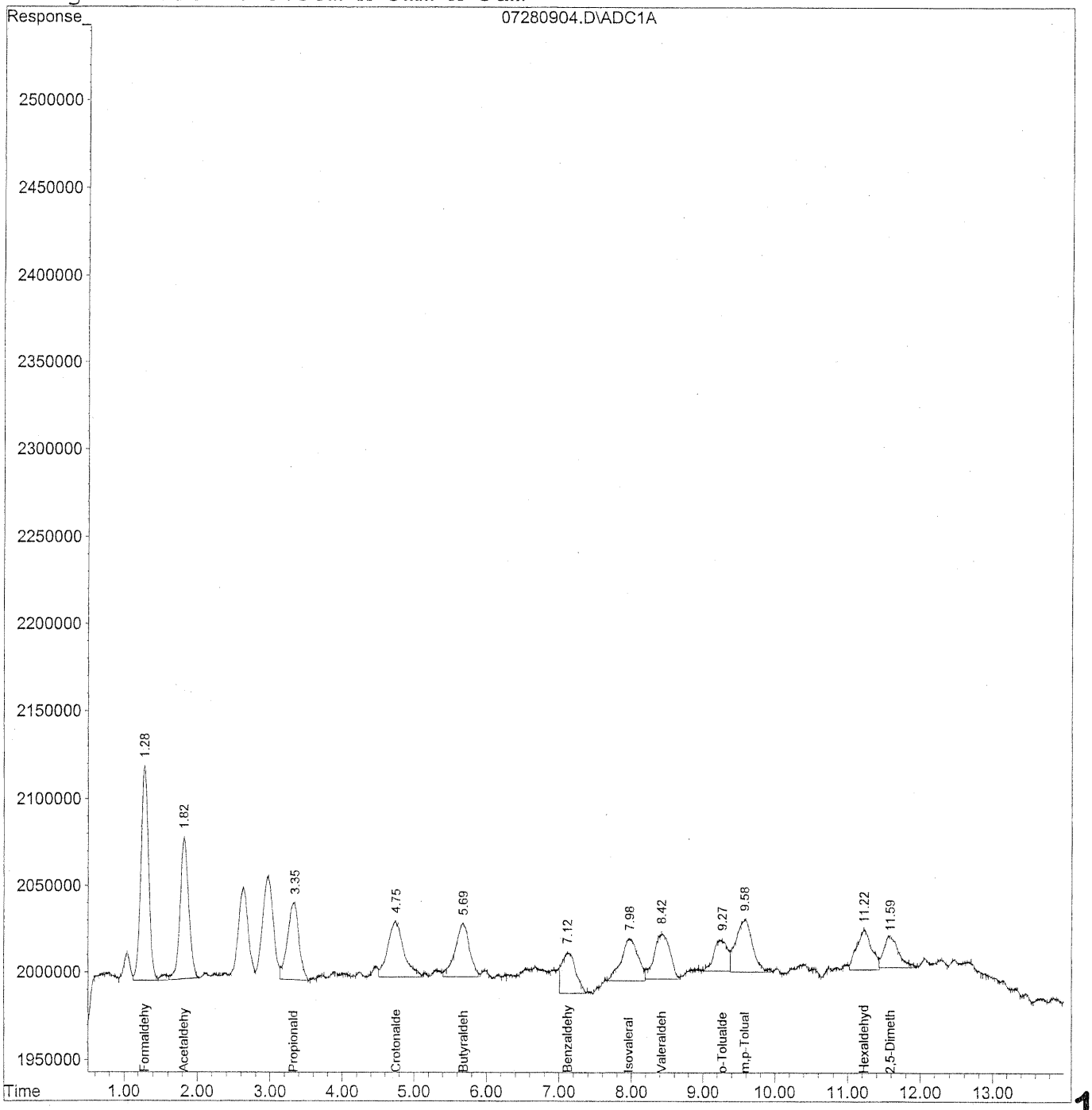
(+) = Expected Retention Time

07280903.D TO110709.M Tue Jul 28 17:30:57 2009

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

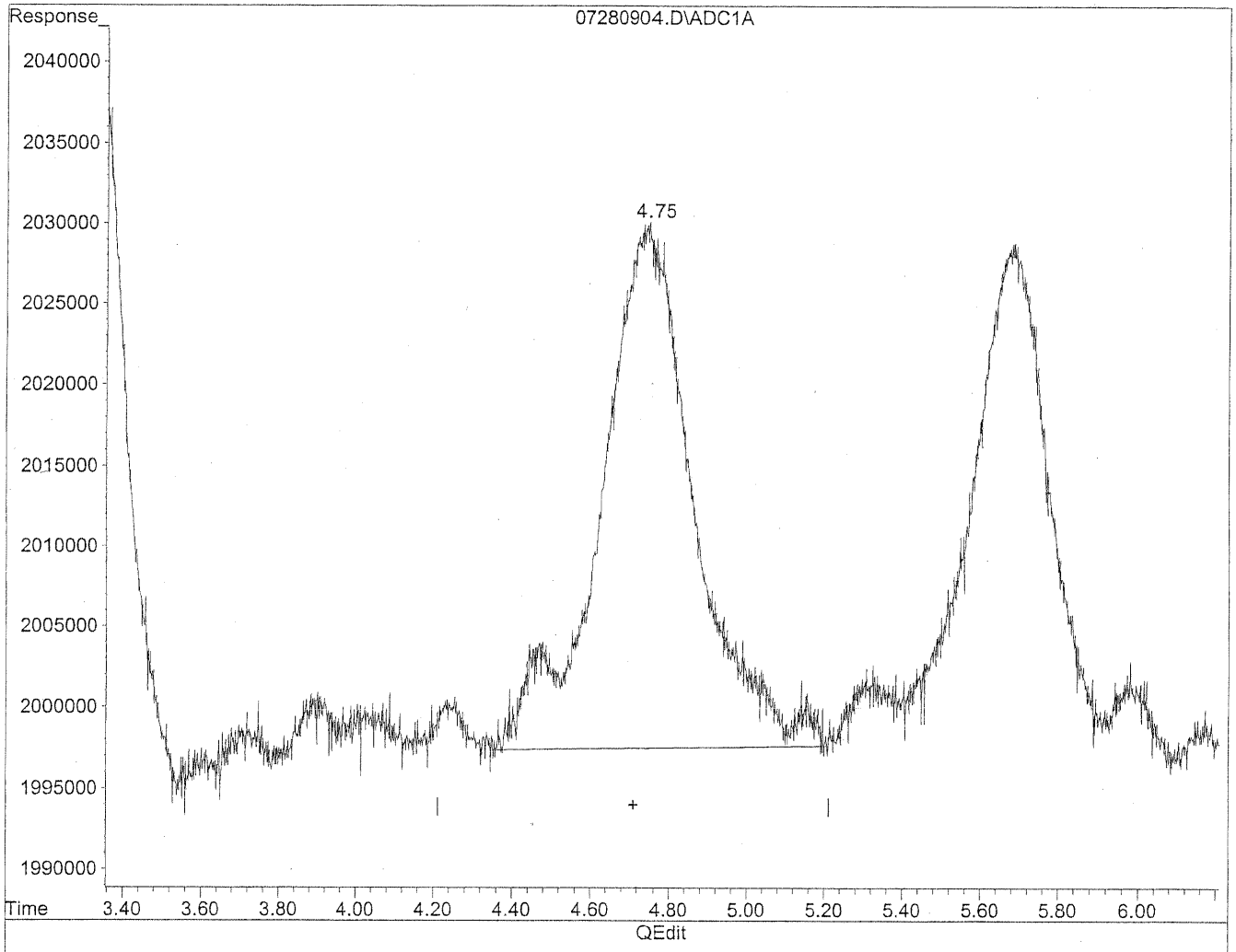
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	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

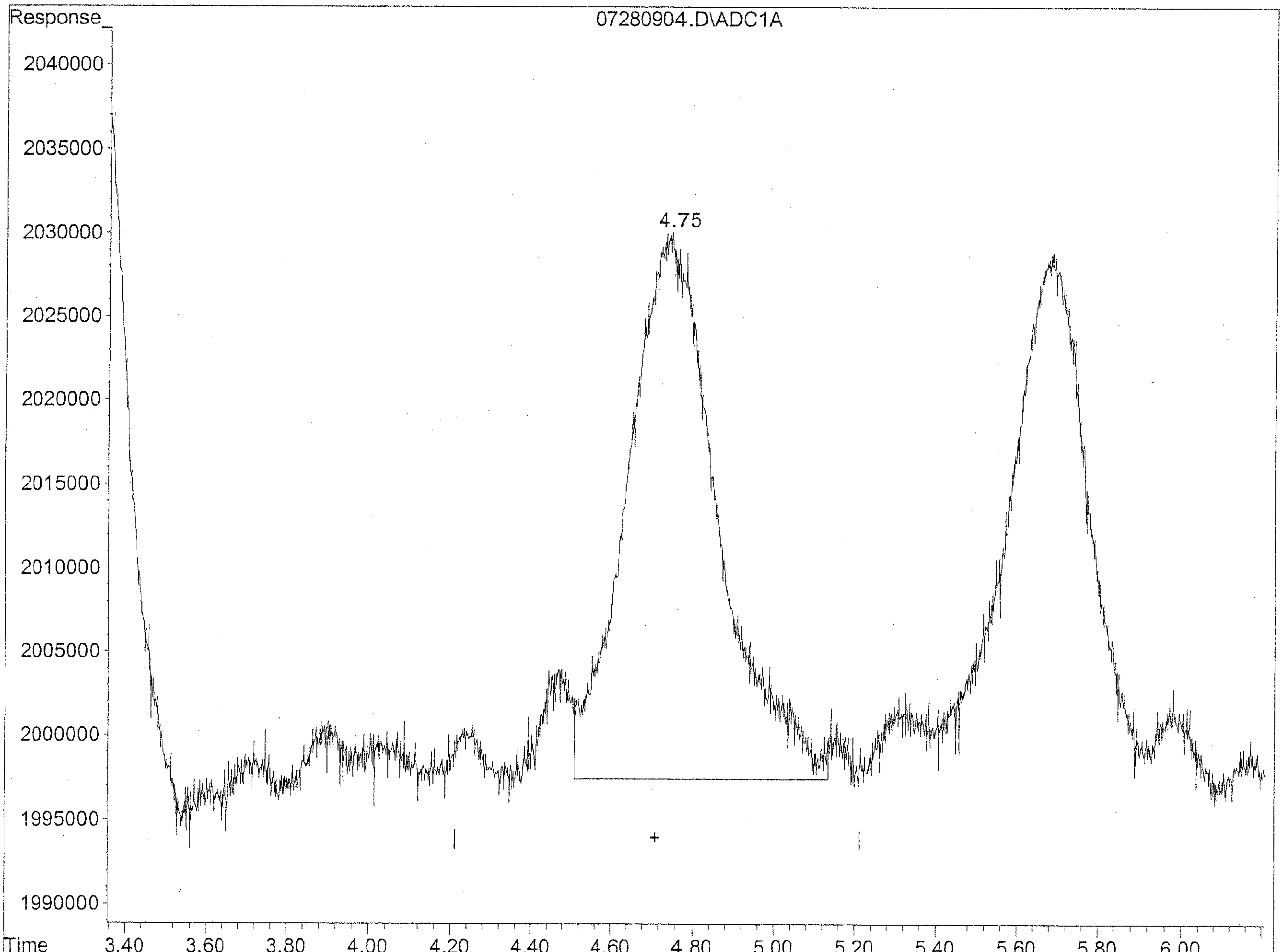
response 5342434

(+) = 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

(4) Crotonaldehyde

4.75min 45.000ng/ml m

response 4974991

HC 7/28/09 C4

KH 7/29/09

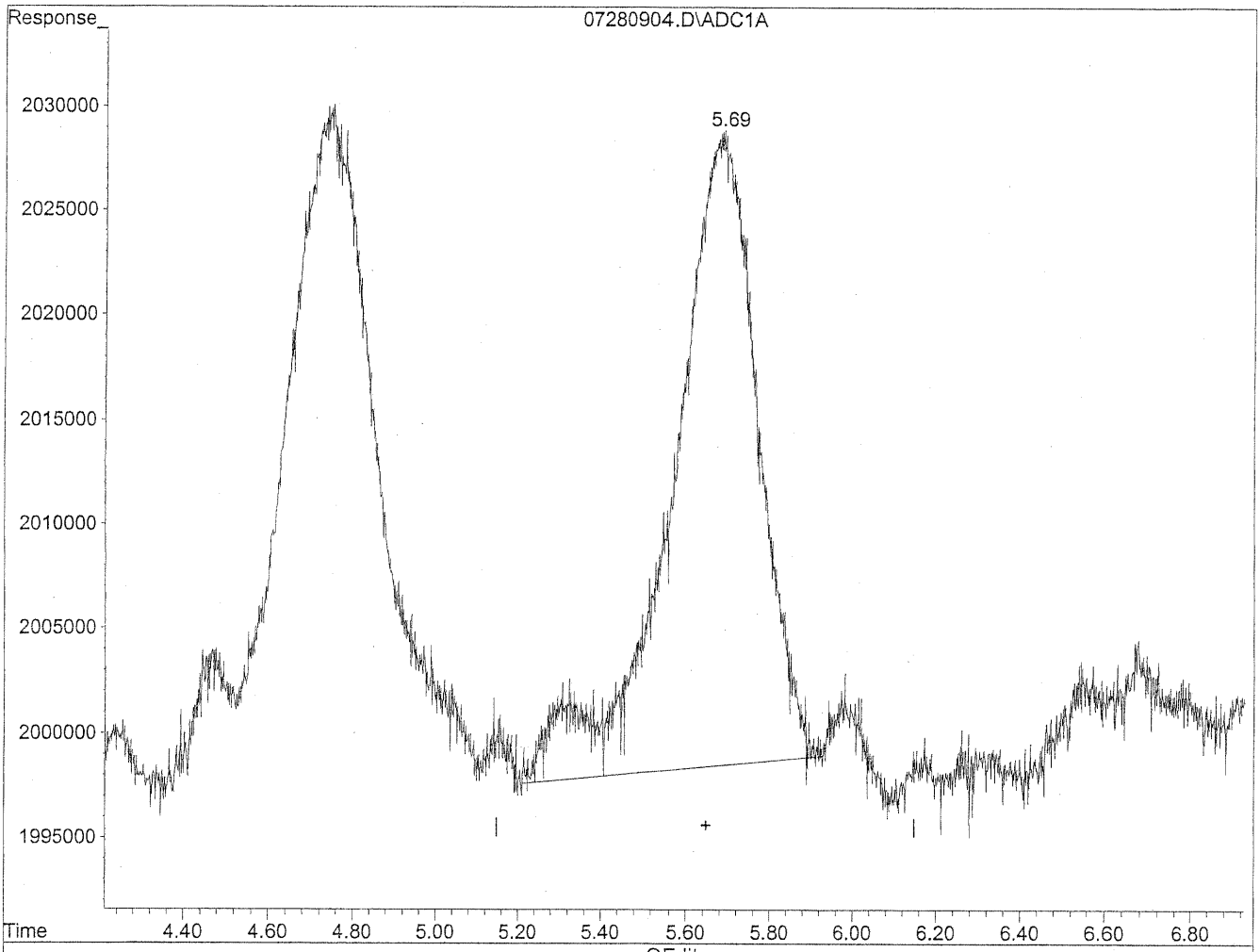
150

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.68min 53.153ng/ml

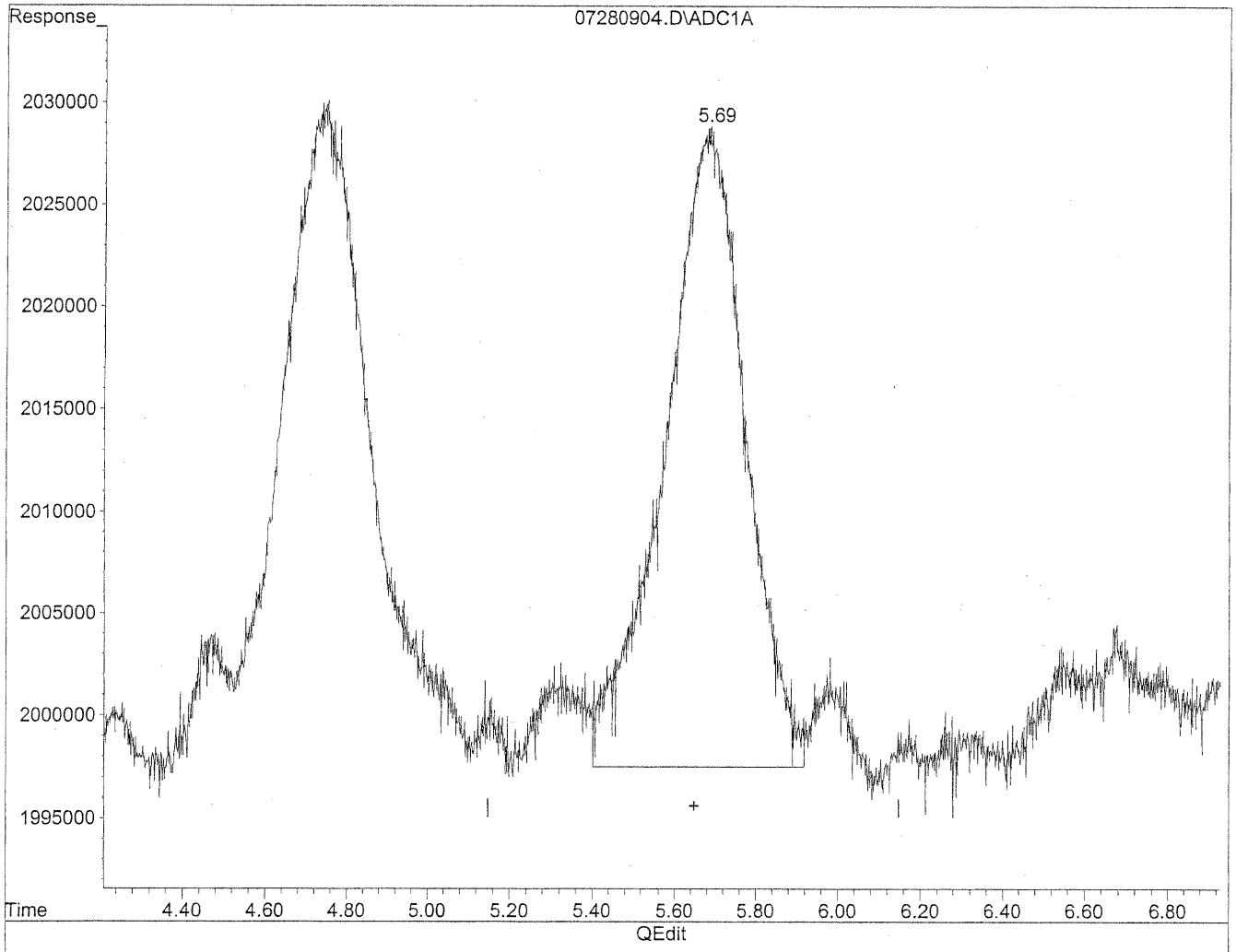
response 4277470

(+) = Expected Retention Time

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

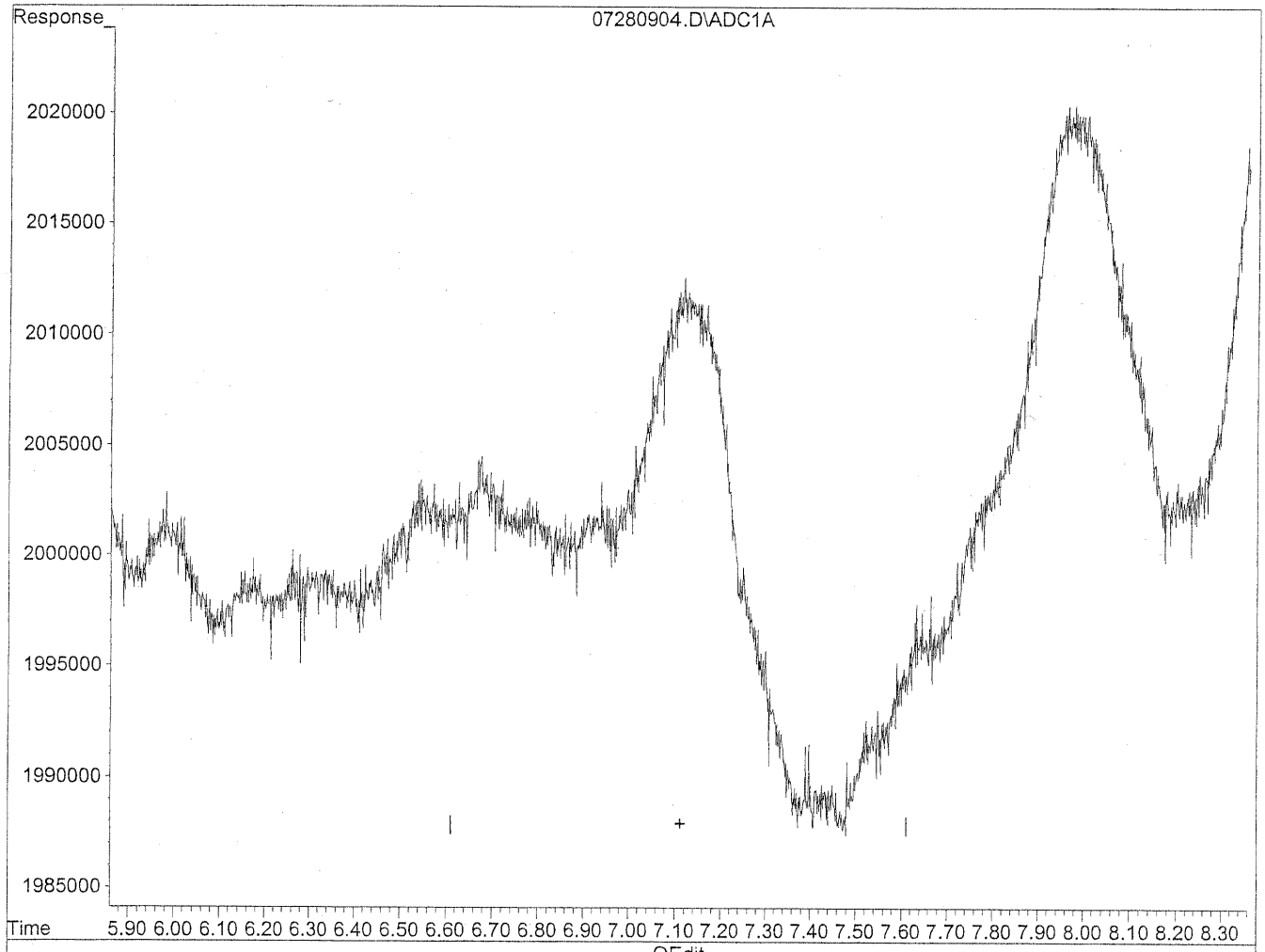
KR 7/29/09

(+) = 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

7.11min 0.000ng/ml

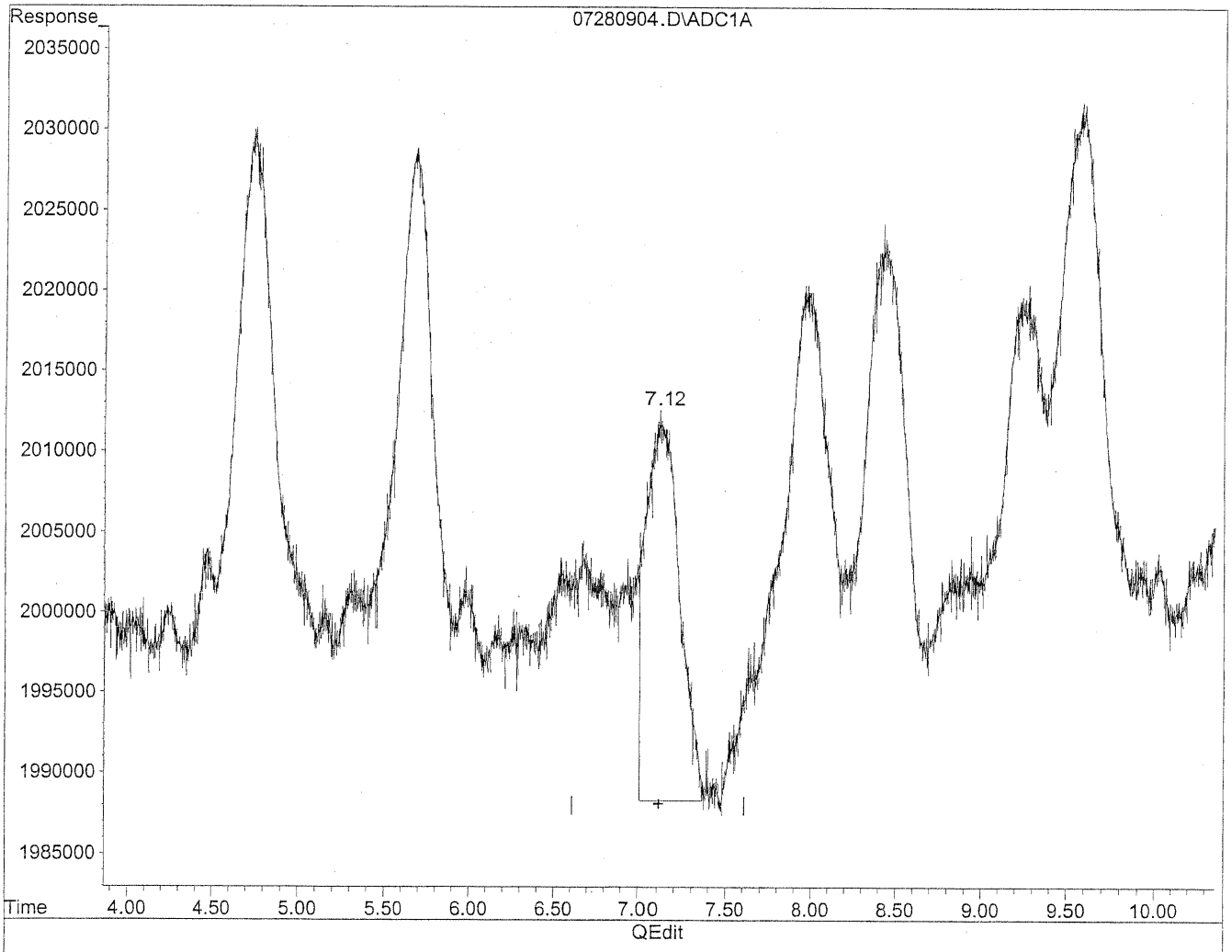
response 0

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

7.12min 48.820ng/ml m

response 3079204

HC 7/28/09 BNL

7/29/09

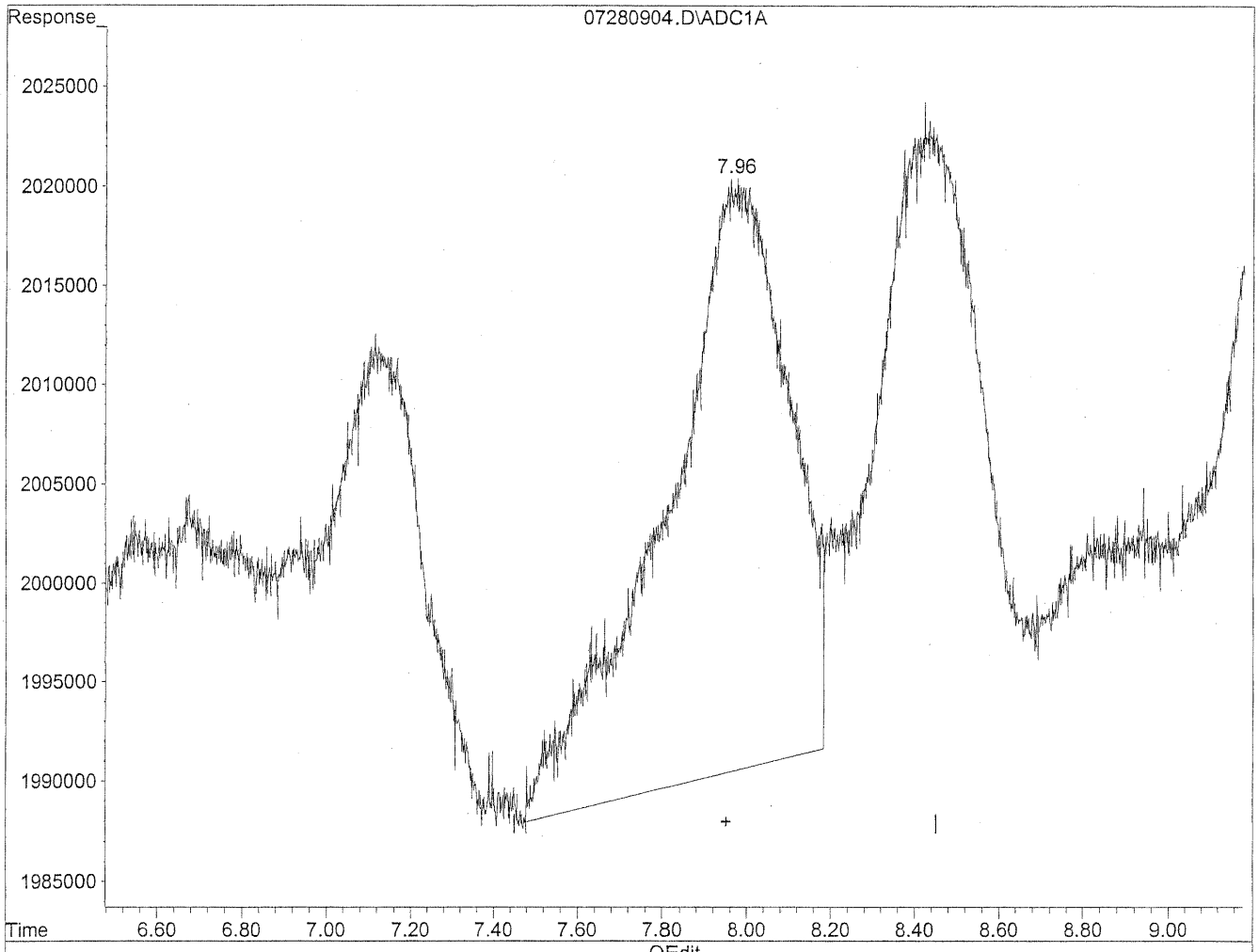
154

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.97min 68.251ng/ml

response 6050534

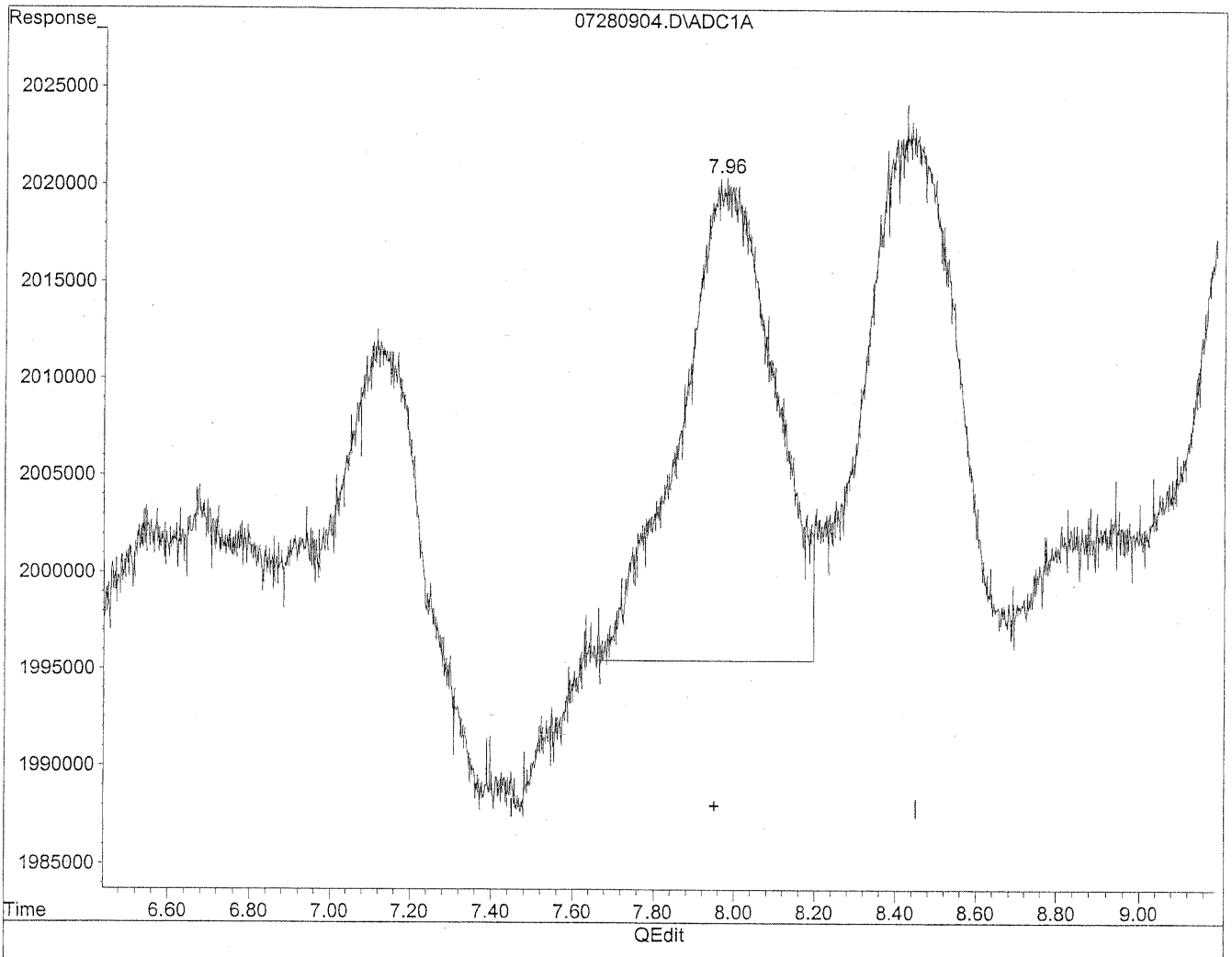
(+) = Expected Retention Time

07280904.D TO110709.M Tue Jul 28 14:24:25 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 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

HC 7/29/09
LC

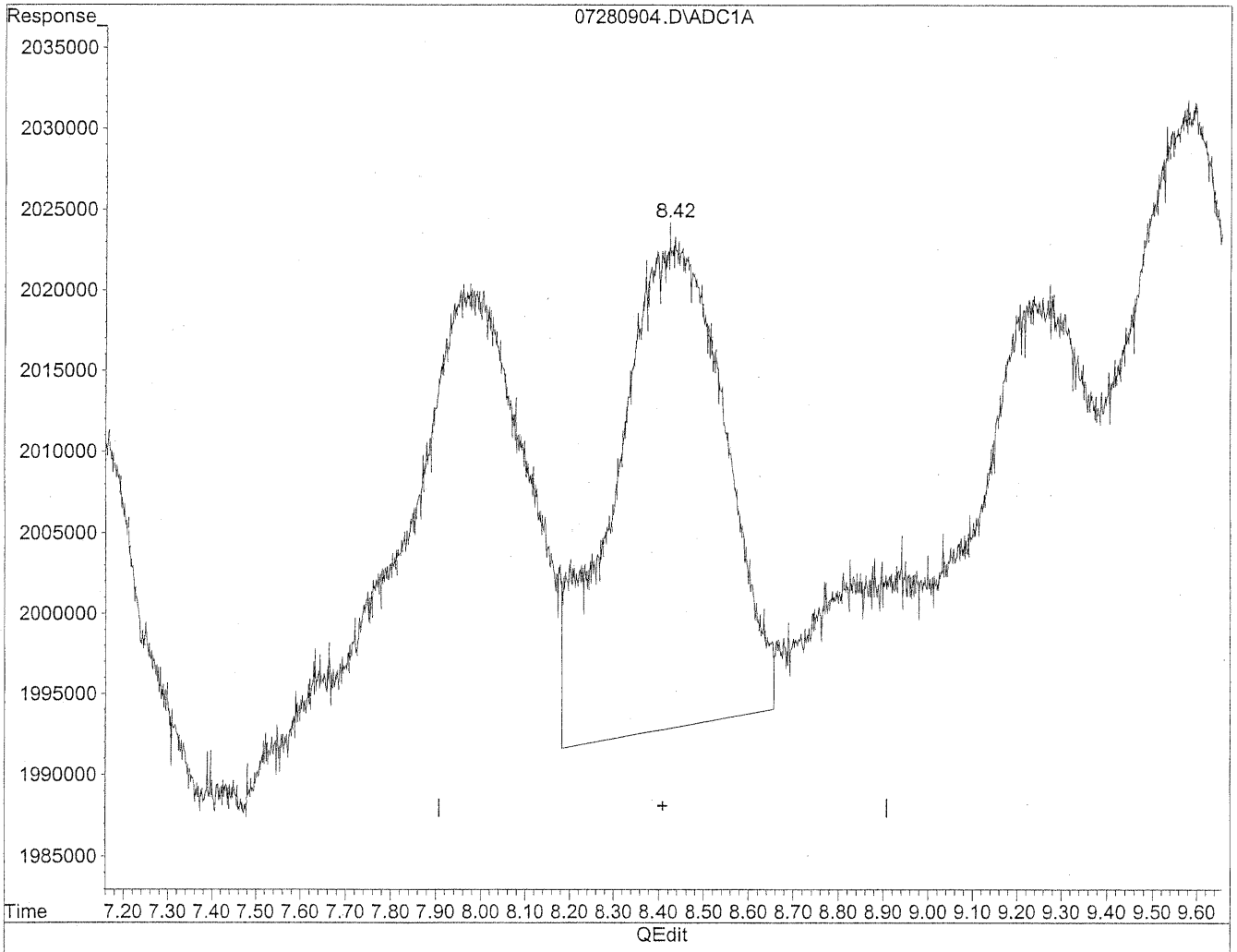
HC 7/30/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

8.43min 61.279ng/ml

response 5091976

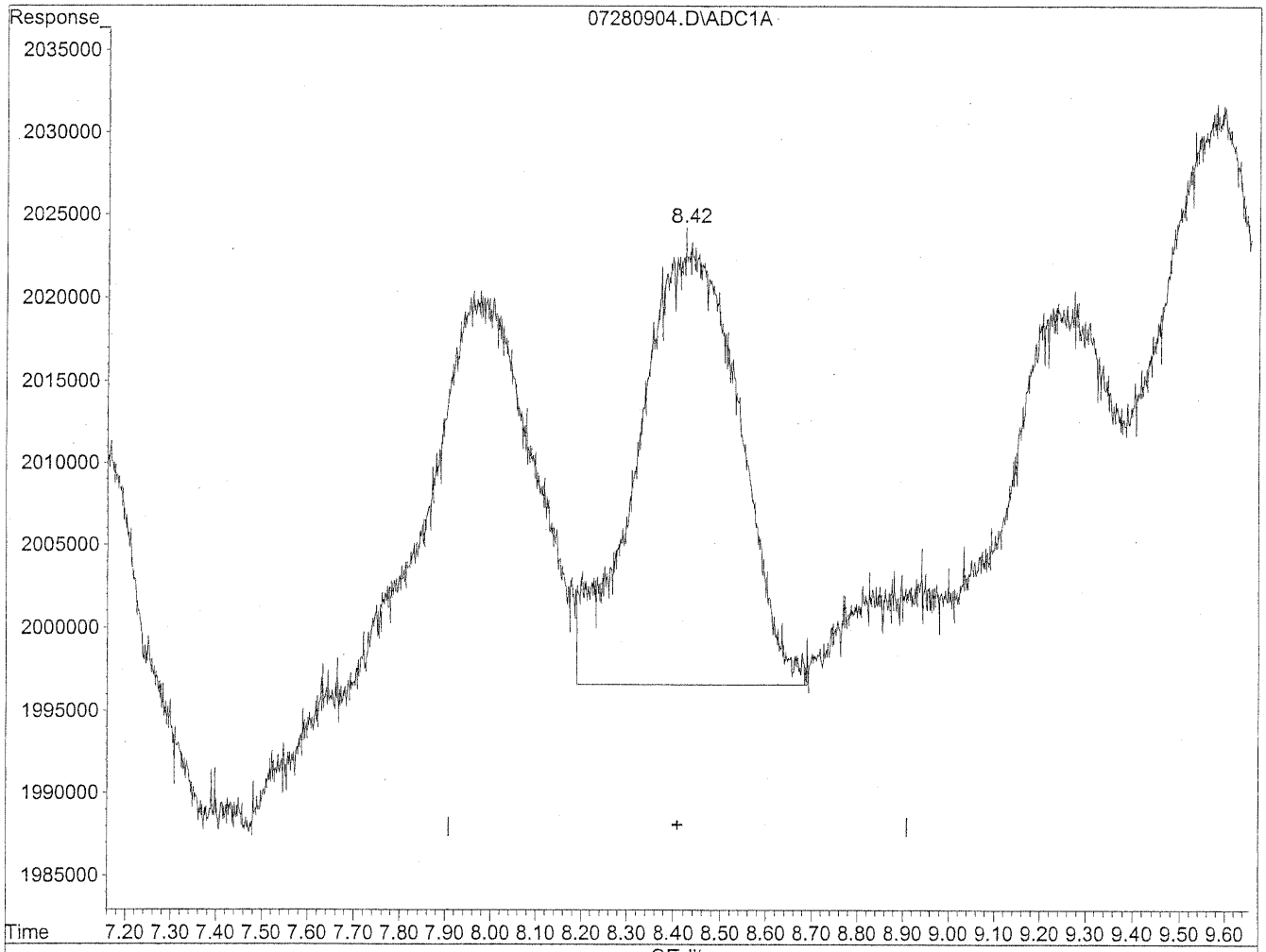
(+) = Expected Retention Time

07280904.D TO110709.M Tue Jul 28 14:24:38 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(8) Valeraldehyde

8.42min 48.445ng/ml m

response 4025564

HC standard LC

KA 7/29/09

158

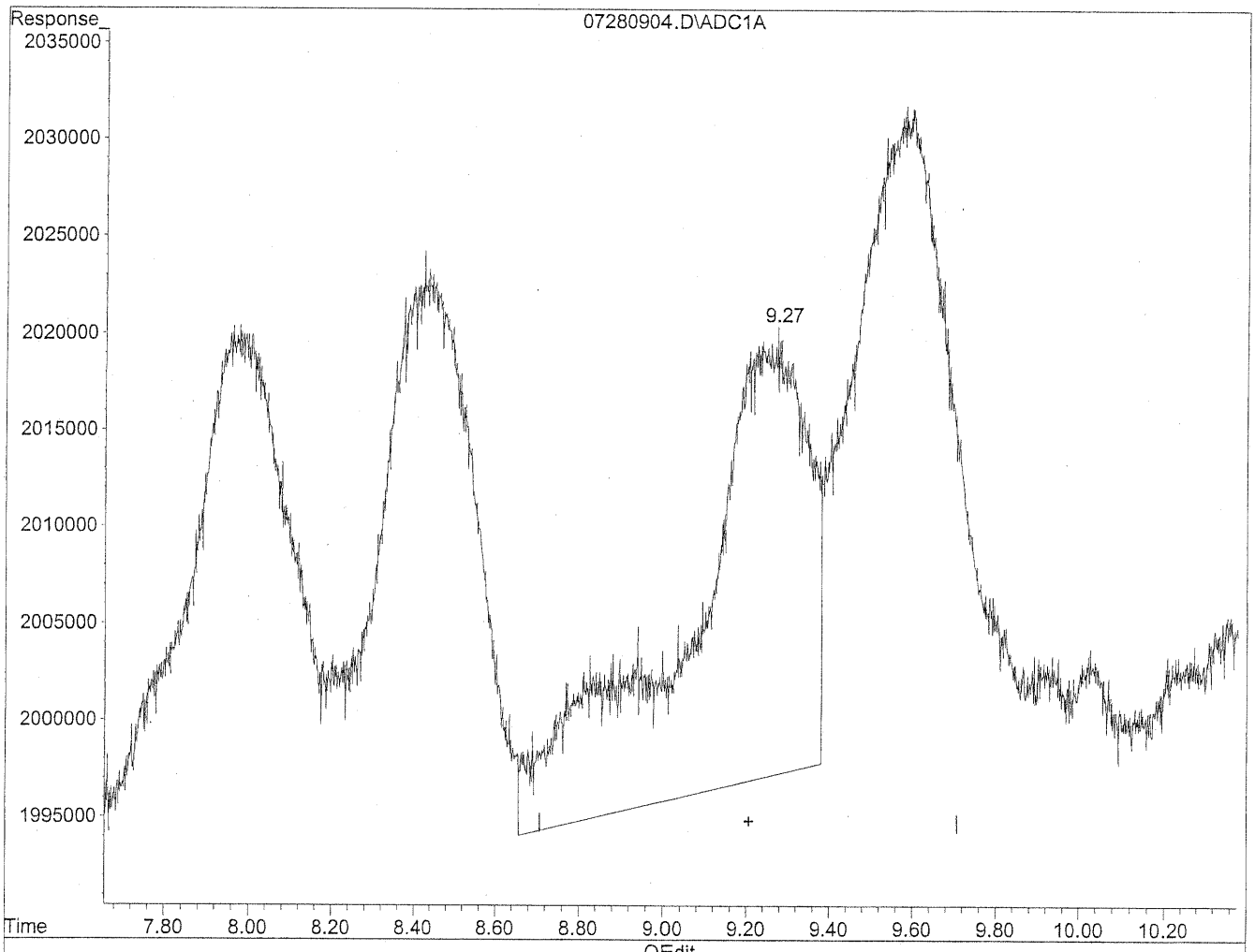
(+) = Expected Retention Time

07280904.D TO110709.M Tue Jul 28 14:24:48 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.24min 84.965ng/ml

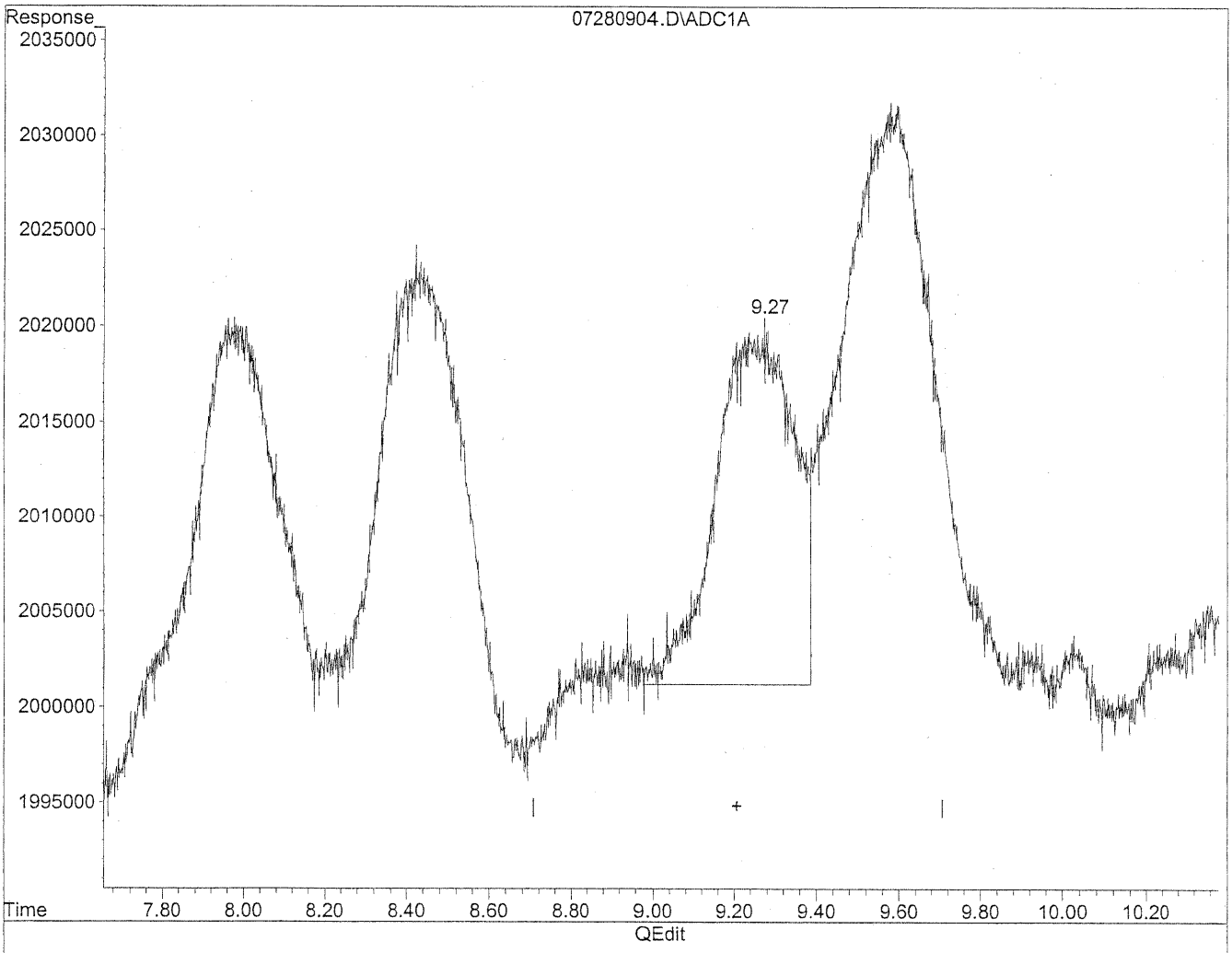
response 4577075

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.27min 45.695ng/ml m

response 2461625

+/C
standard
LC

KP 7/29/09

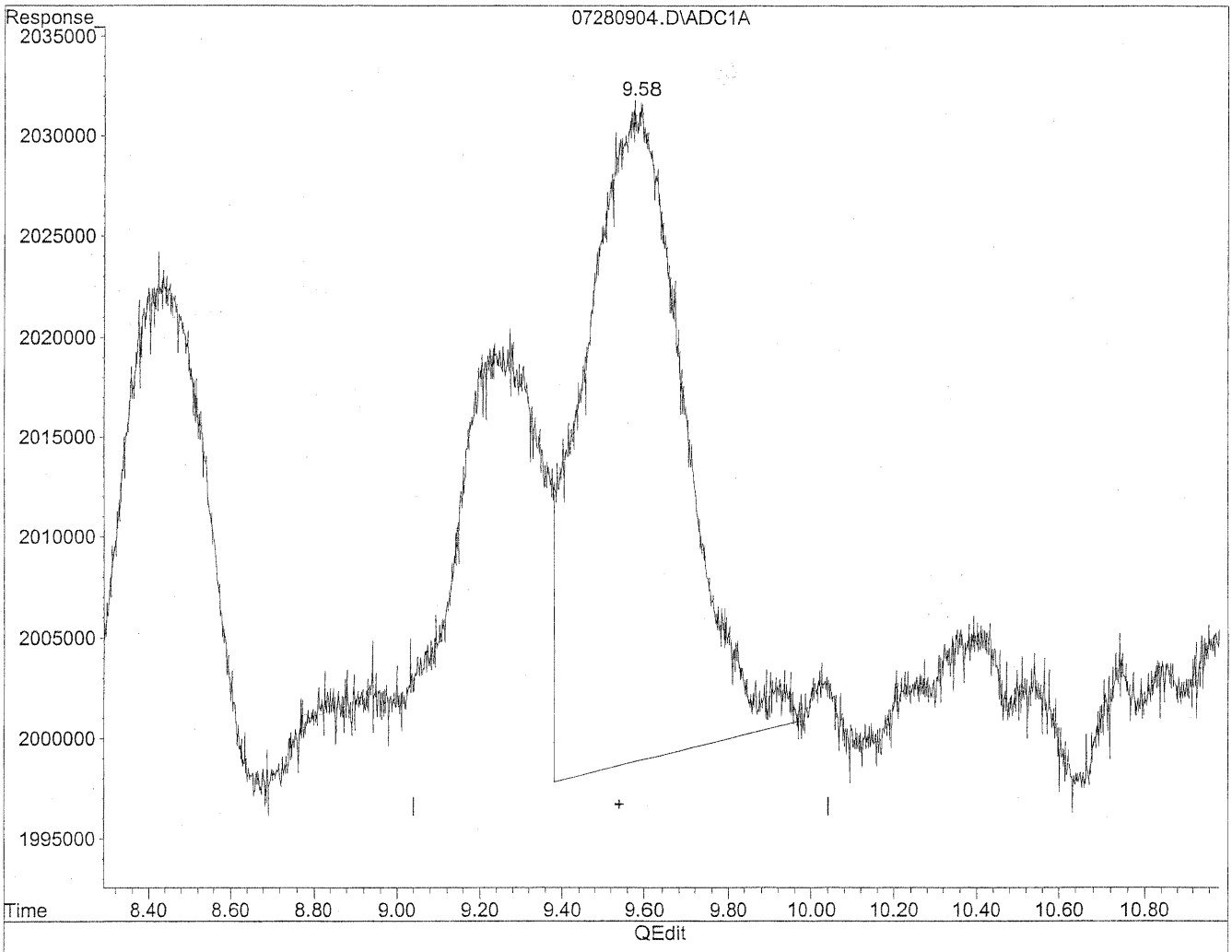
160

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

9.59min 100.987ng/ml

response 5439618

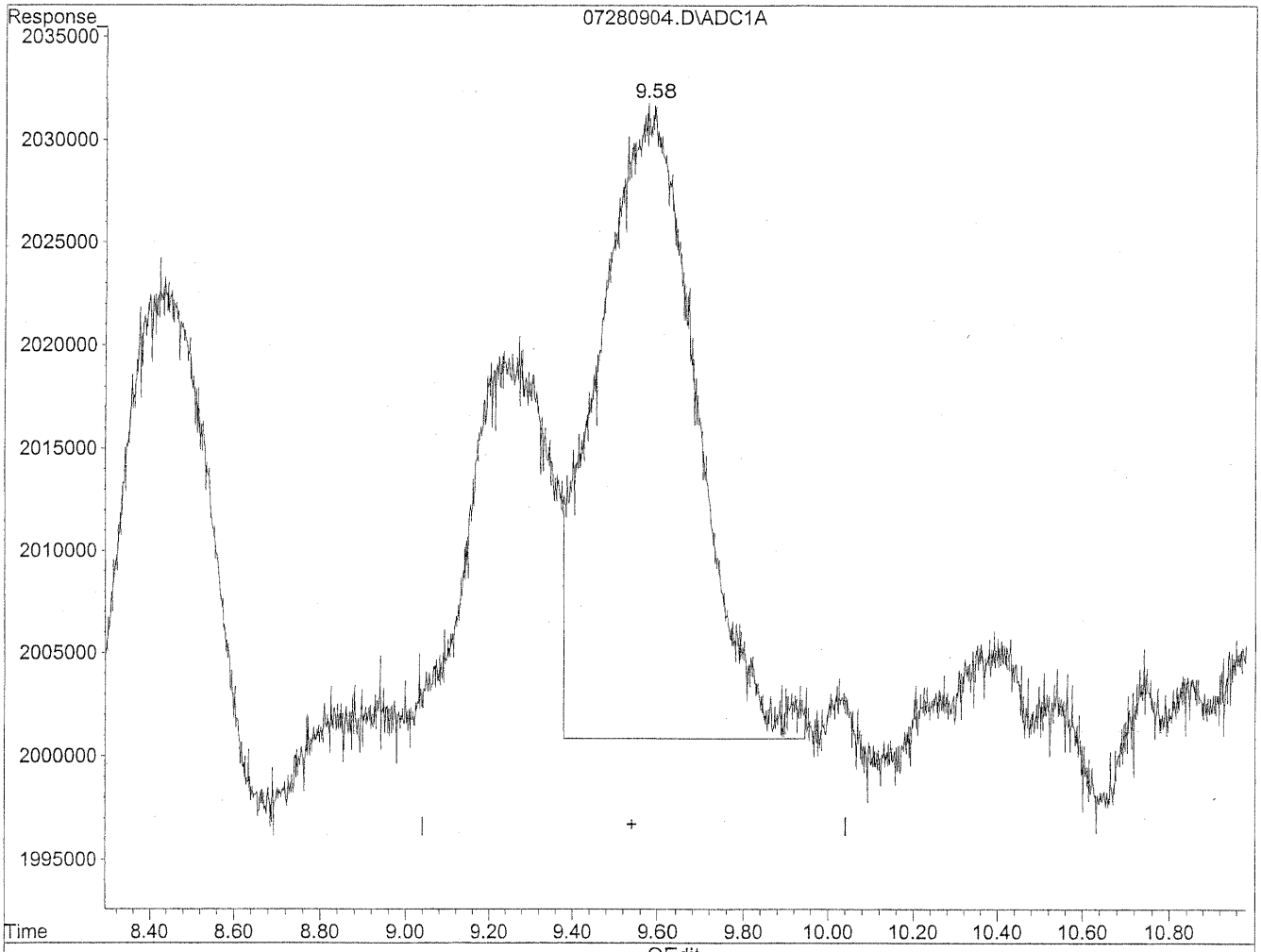
161

(+) = 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

(10) m,p-Tolualdehyde

9.58min 90.915ng/ml m

response 4897087

HC 7/28/09

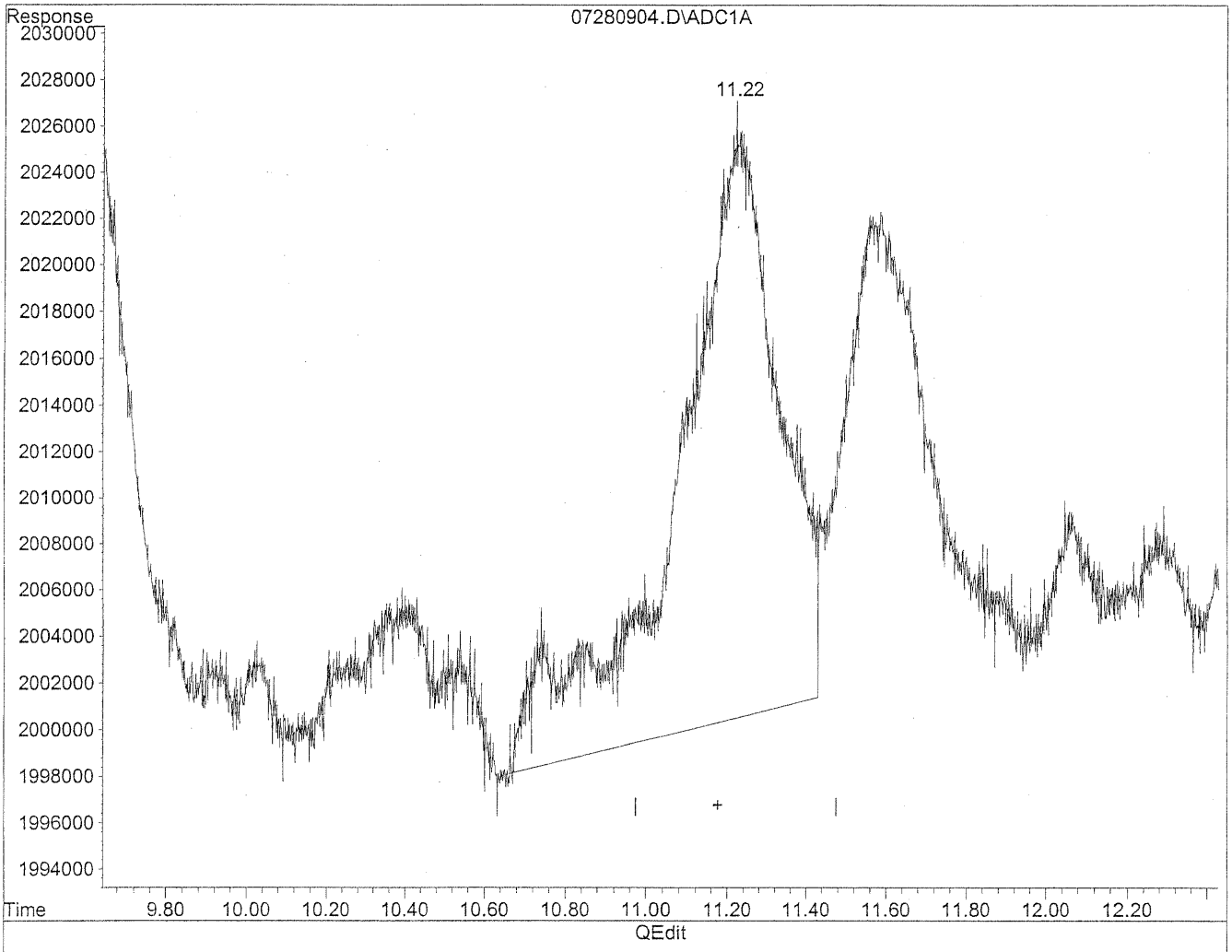
KR 7/29/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

11.23min 66.912ng/ml

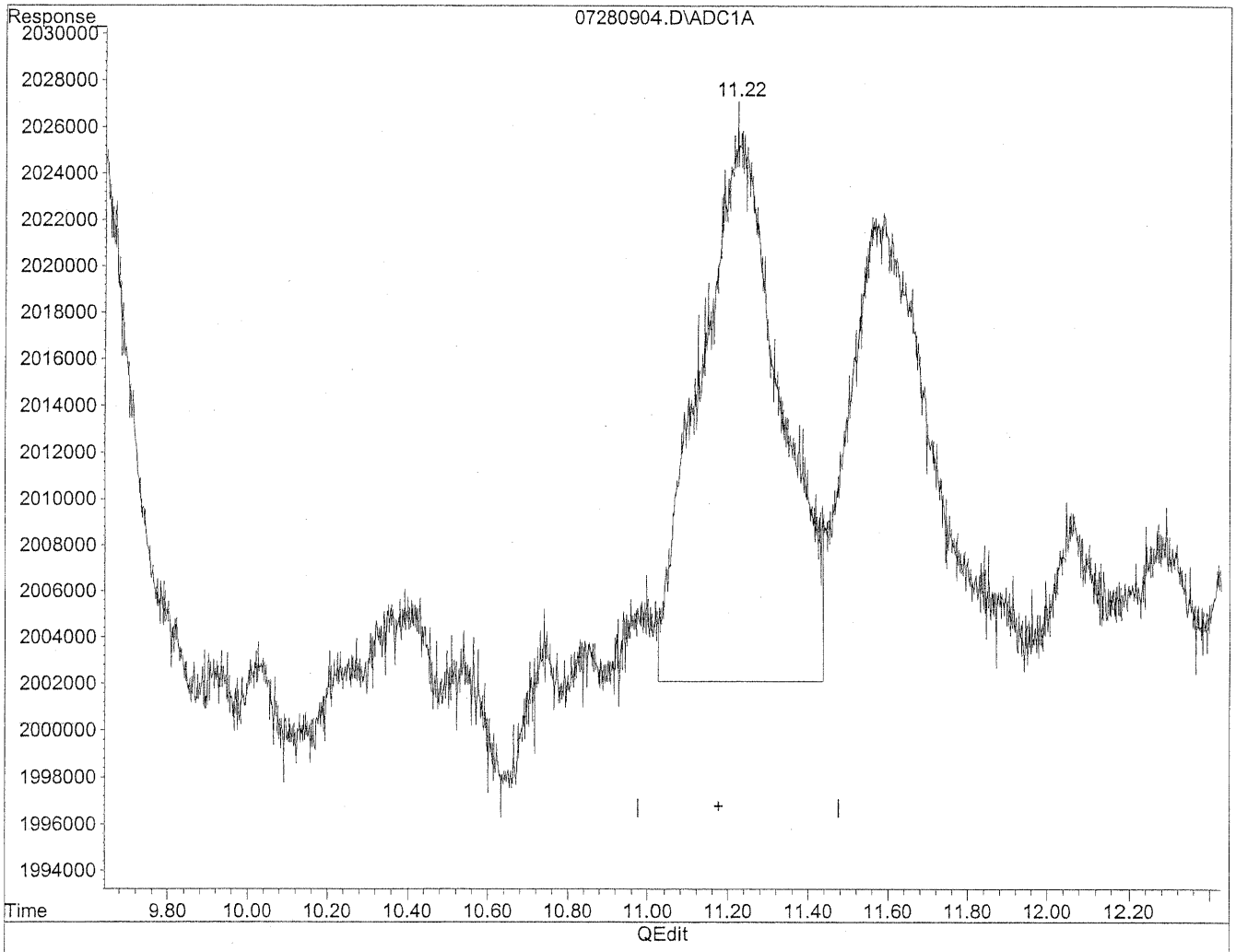
response 4492347

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

11.22min 49.079ng/ml m

response 3295067

HC 7/28/09 SH

KR 7/29/09

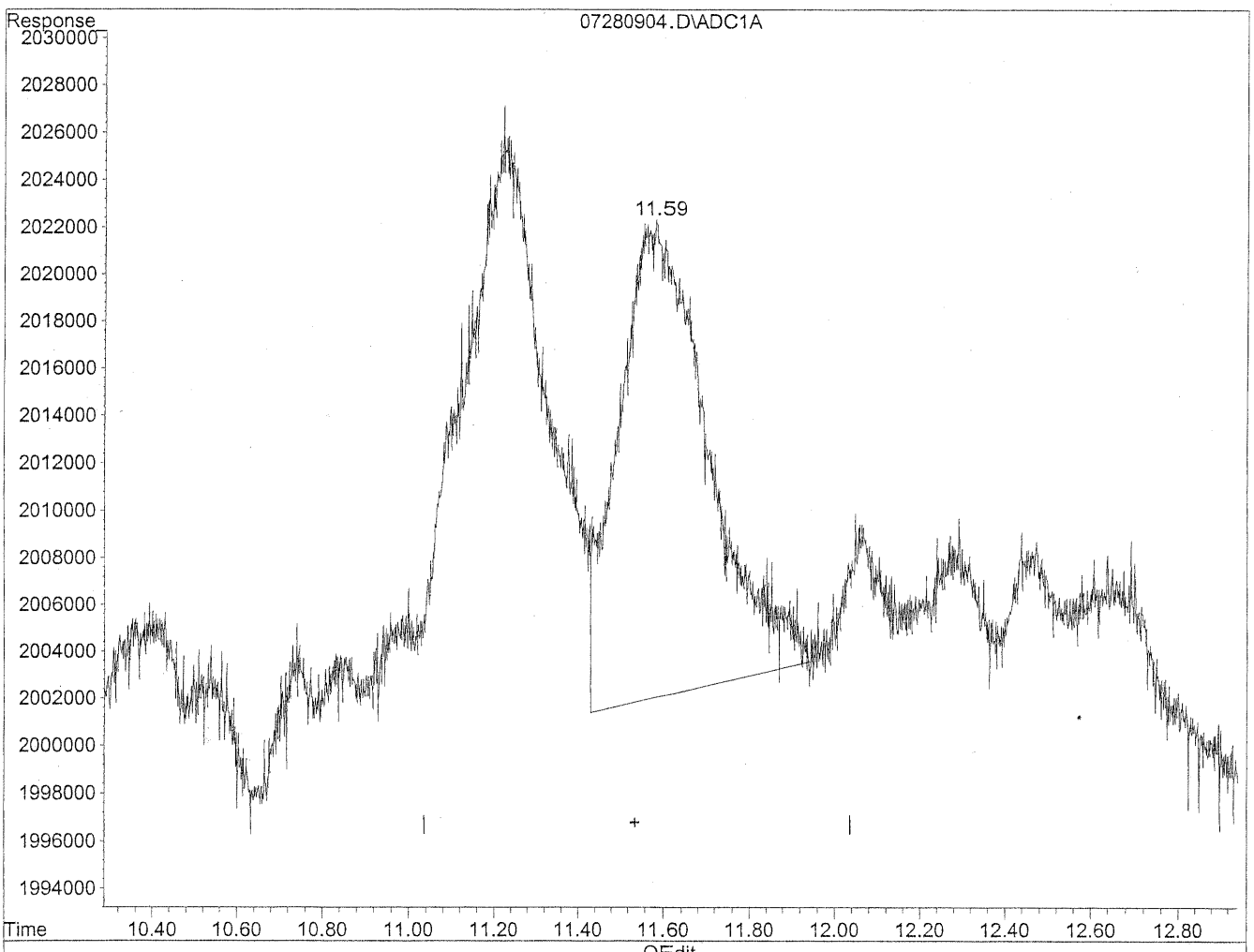
(+) = Expected Retention Time

07280904.D TO110709.M Tue Jul 28 14:25:22 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



QEdit

(12) 2,5-Dimethylbenzaldehyde

11.58min 55.789ng/ml

response 2897339

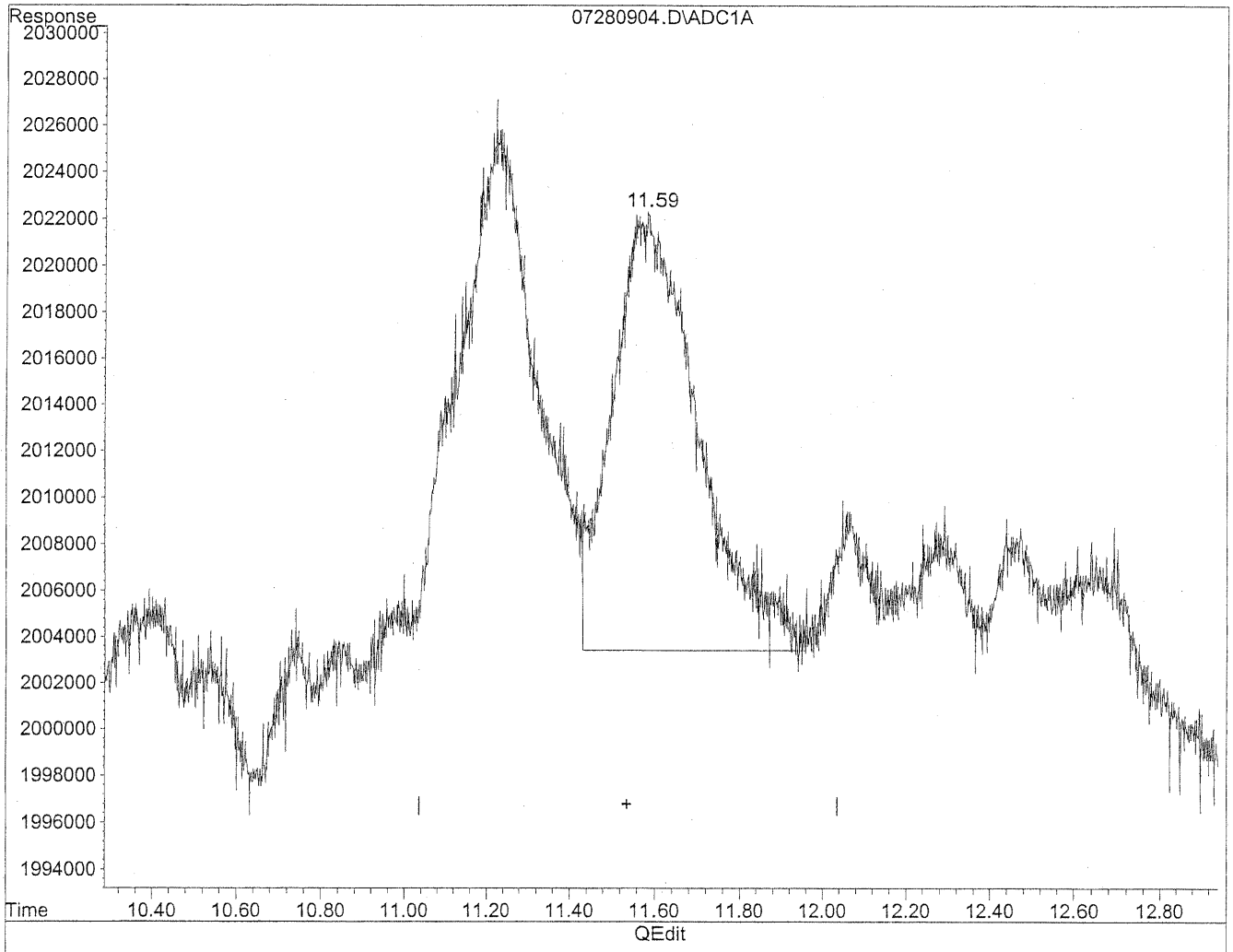
(+) = Expected Retention Time

07280904.D TO110709.M Tue Jul 28 14:25:29 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280904.D Vial: 4
Acq On : 28 Jul 2009 9:24 am Operator: HC
Sample : 50ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:23 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

11.59min 50.169ng/ml m

response 2605446

HC 9/28/09 LC

KF 7/29/09

166

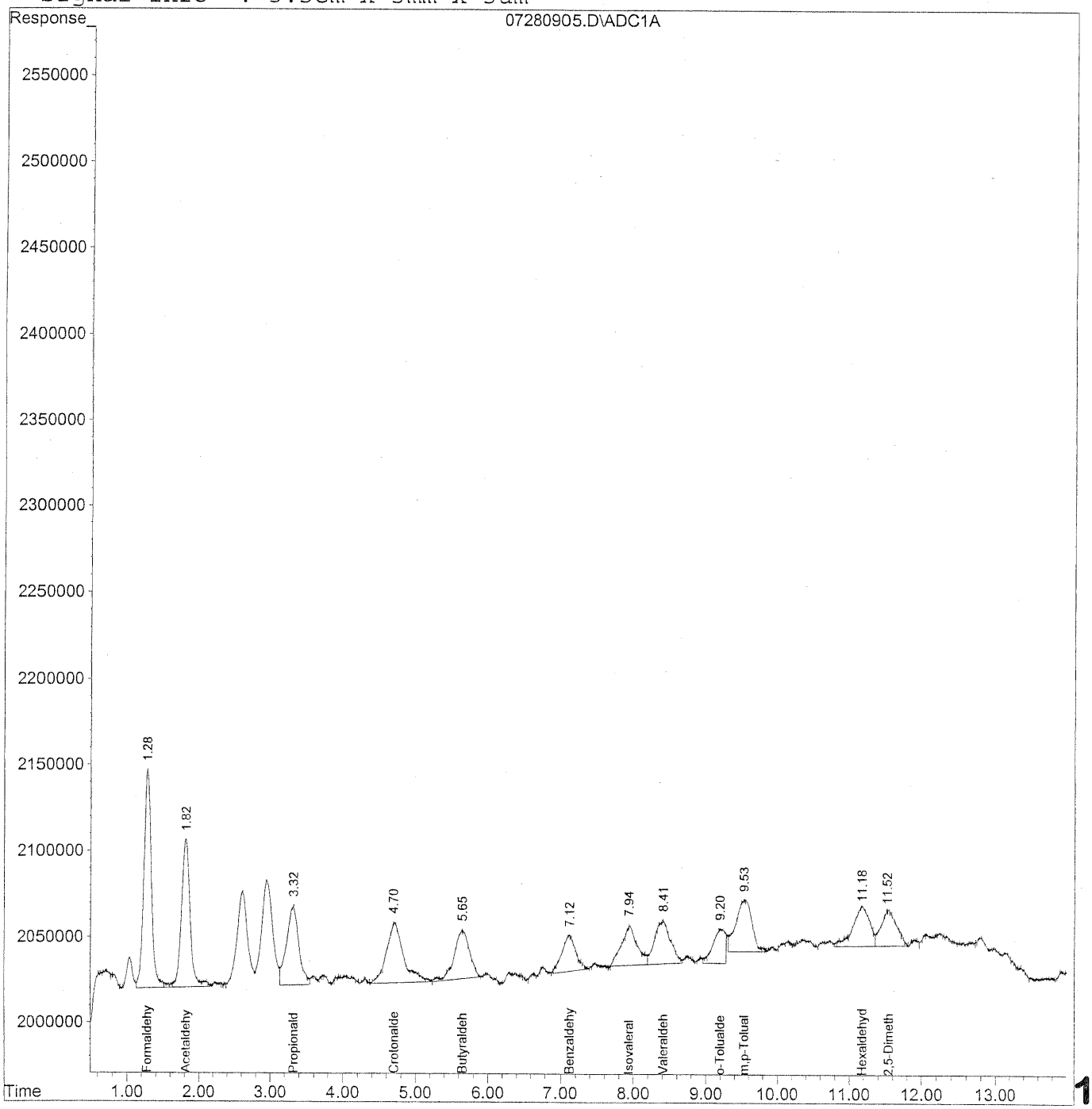
(+) = Expected Retention Time
07280904.D TO110709.M Tue Jul 28 14:25:43 2009

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



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

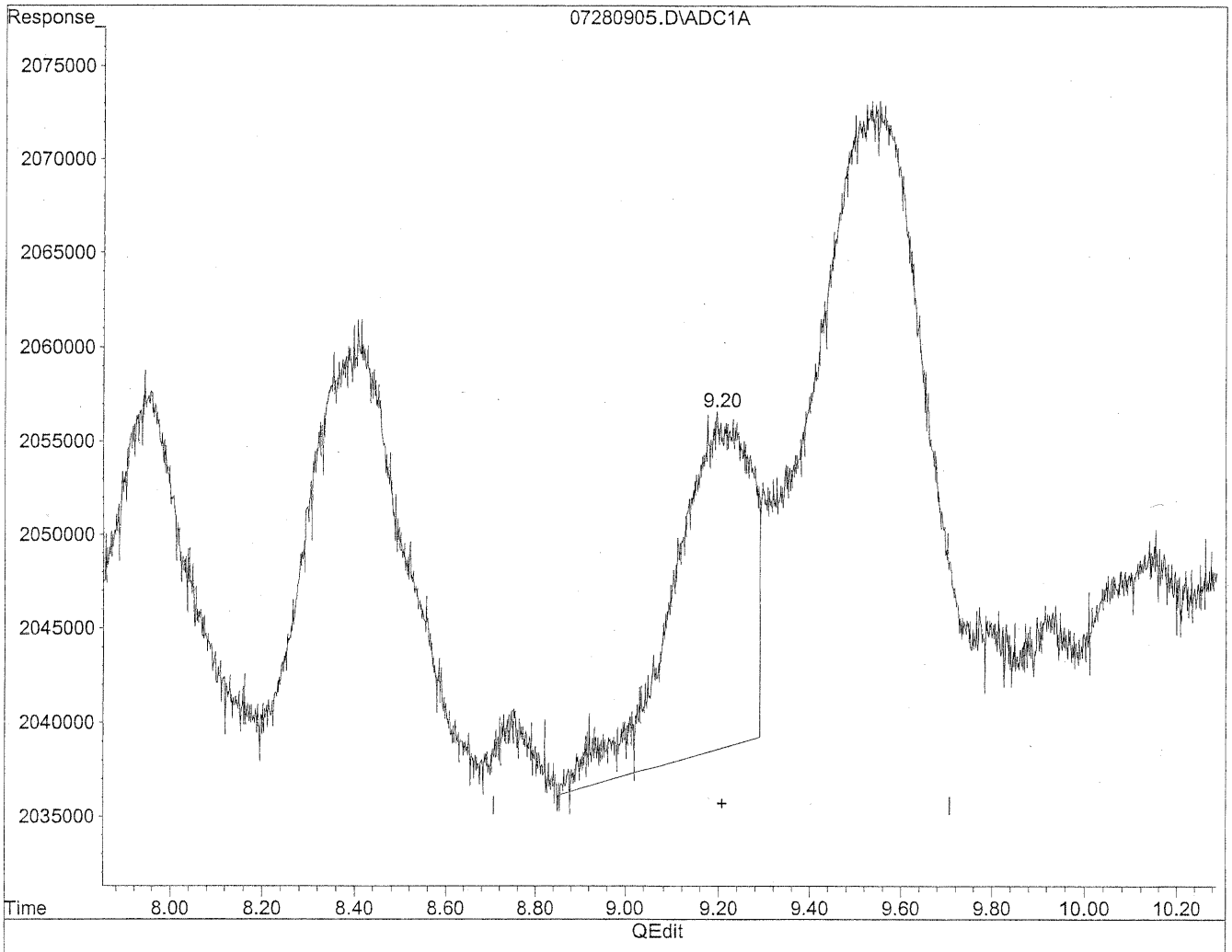
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.28	9305088	52.992	ng/ml
2)	Acetaldehyde	1.81	7389770	54.780	ng/ml
3)	Propionaldehyde	3.31	5442713	52.947	ng/ml
4)	Crotonaldehyde	4.71	5754474	52.051	ng/ml
5)	Butyraldehyde	5.65	4119144	51.185	ng/ml
6)	Benzaldehyde	7.11	2732056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

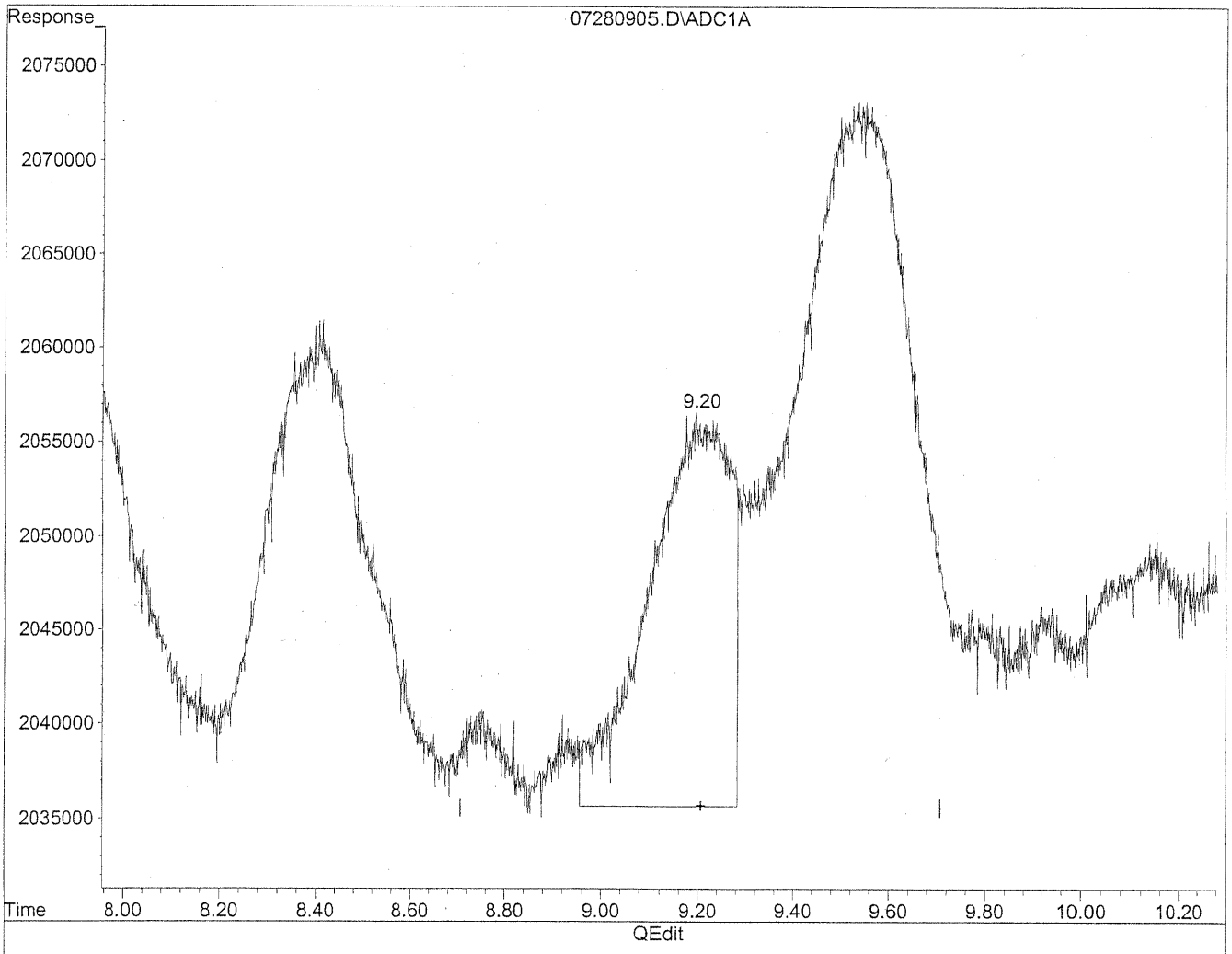
response 2078690

(+) = 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

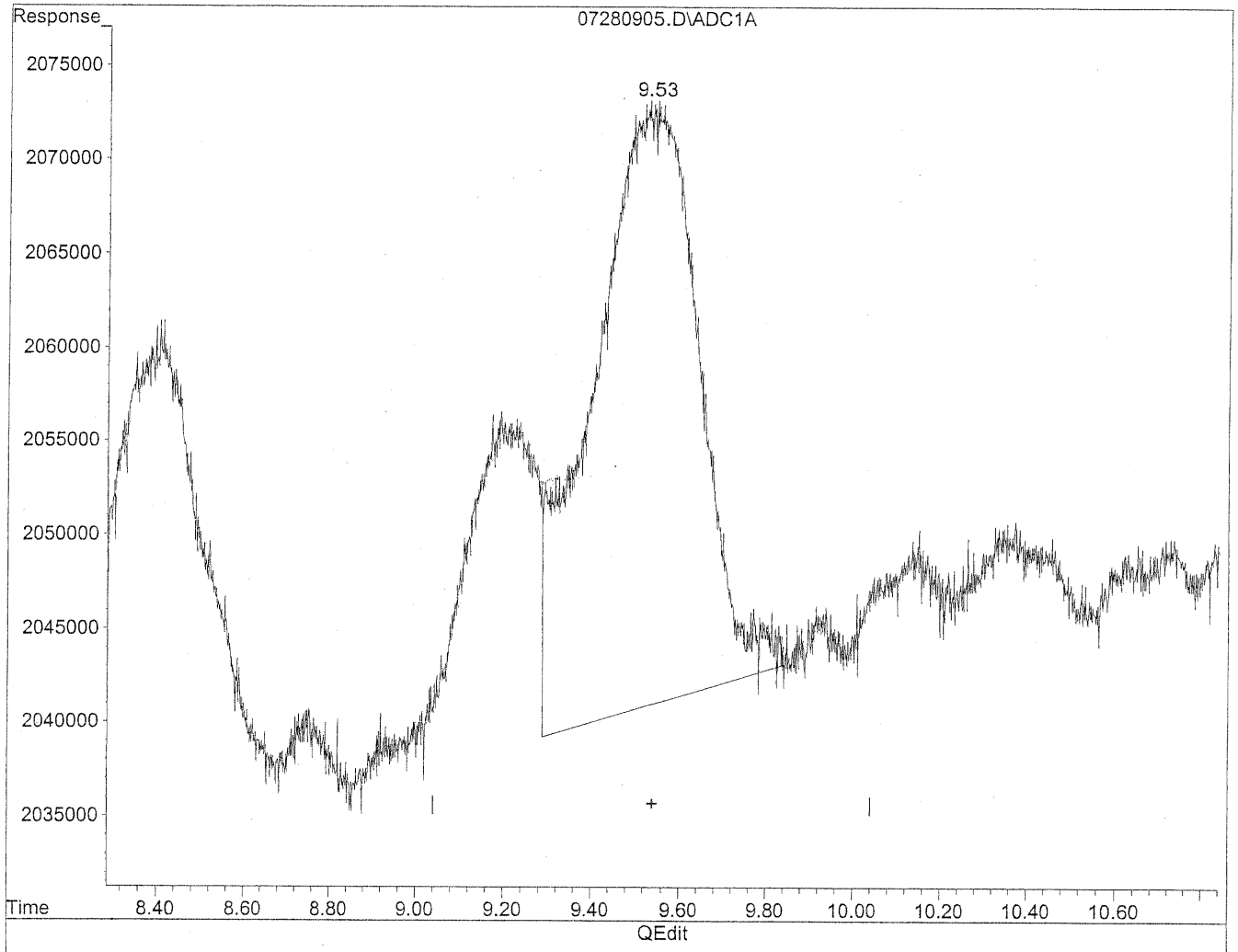
(+) = Expected Retention Time

07280905.D TO110709.M Tue Jul 28 17:33:06 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 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

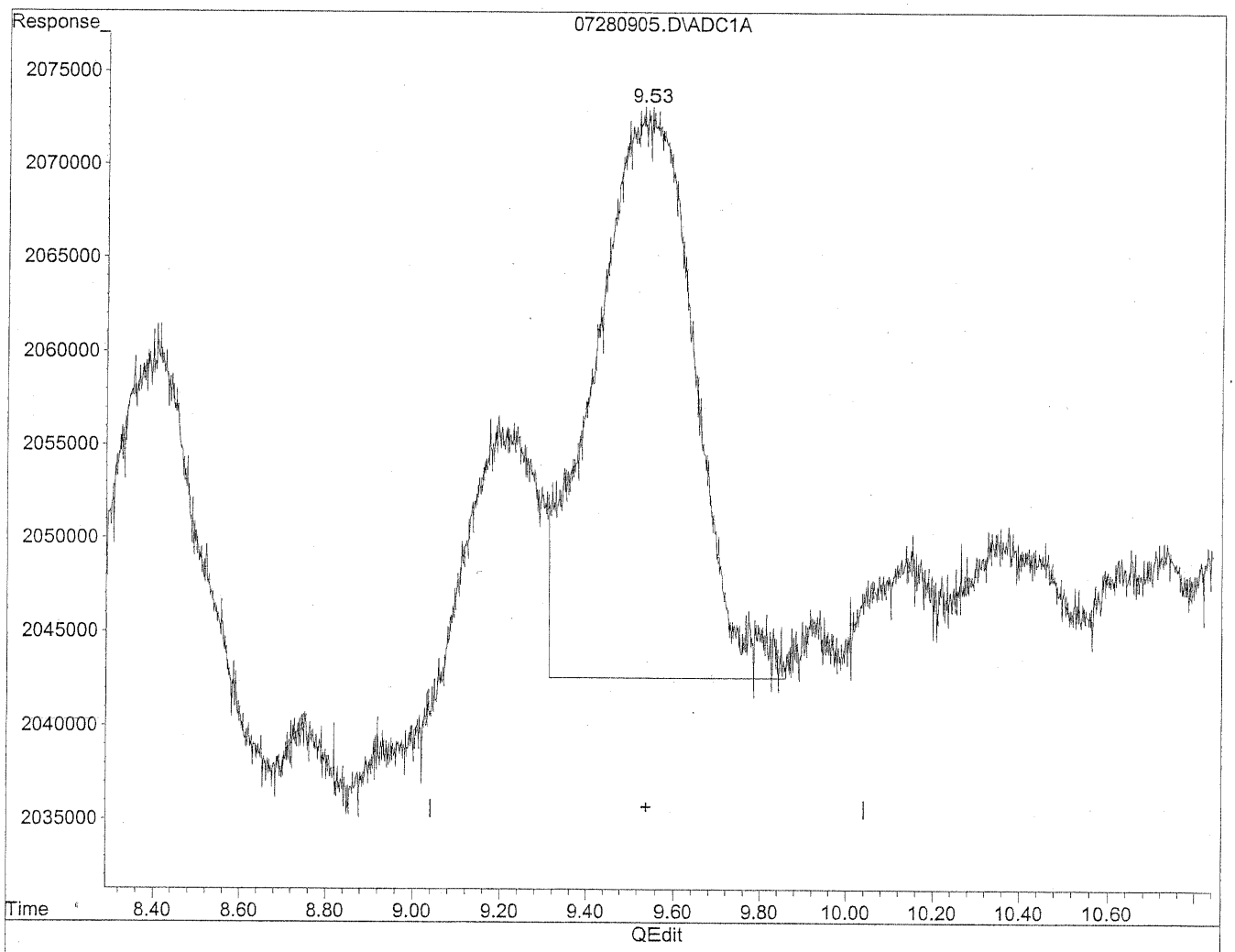
171

(+) = 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

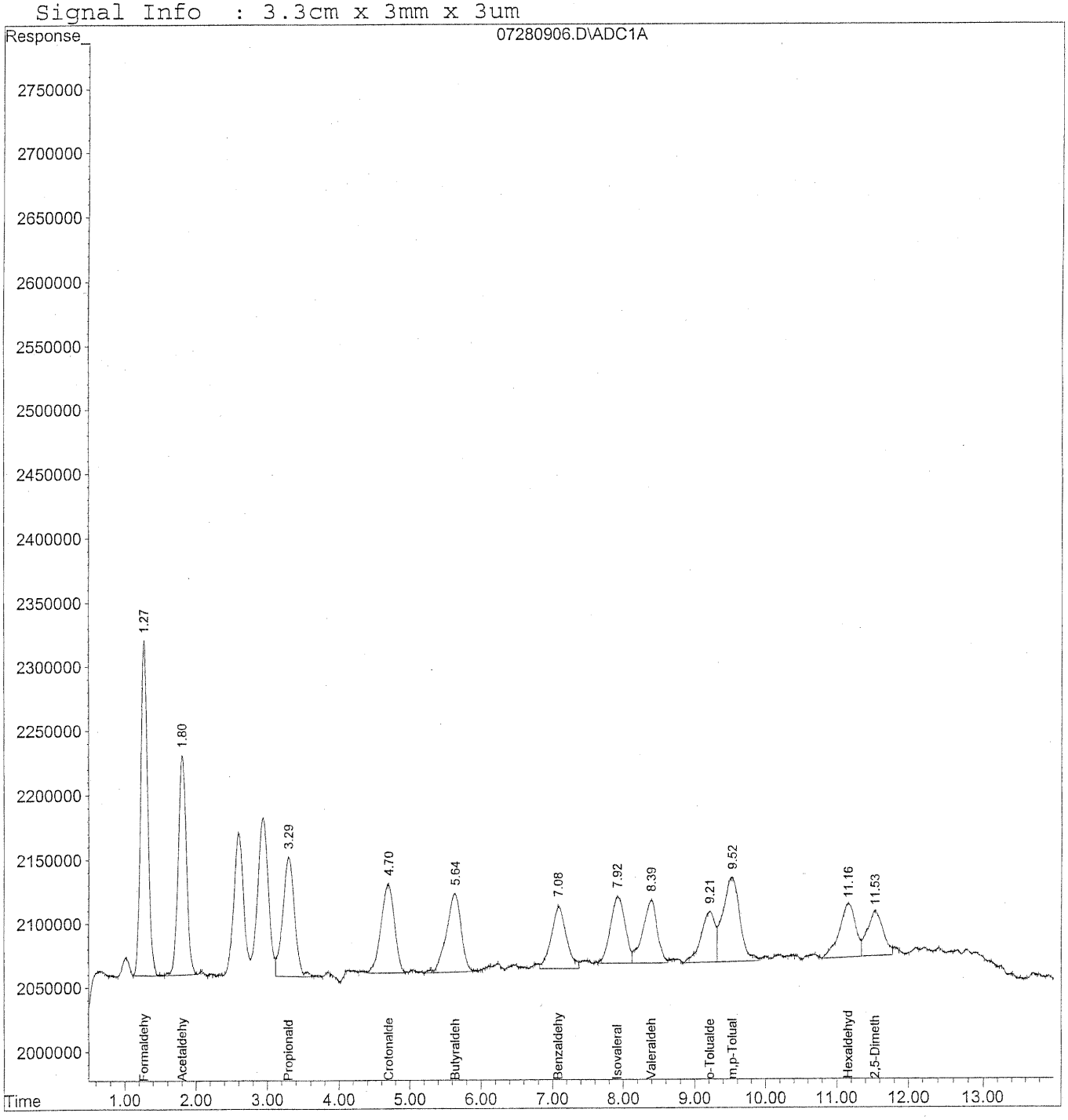
7/29/09

(+) = 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: 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



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280906.D Vial: 6
Acq On : 28 Jul 2009 9:54 am Operator: HC
Sample : 100ng/ml TO11A Std S21-07270905 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: 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

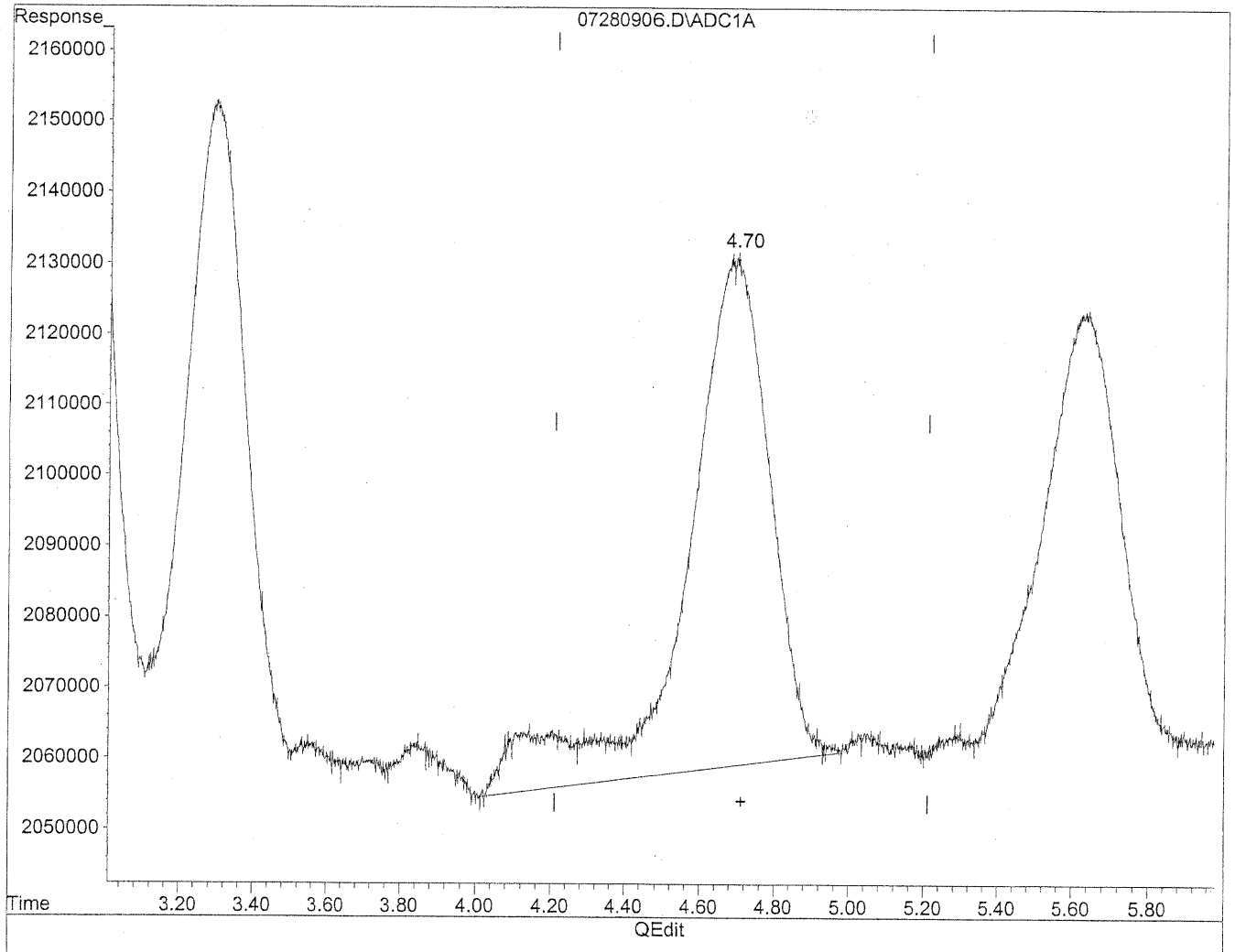
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.27	18283557	104.125	ng/ml
2)	Acetaldehyde	1.80	13784712	102.185	ng/ml
3)	Propionaldehyde	3.29	10870707	105.751	ng/ml
4)	Crotonaldehyde	4.70	9346475	84.541	ng/mlm
5)	Butyraldehyde	5.63	8839595	109.842	ng/ml
6)	Benzaldehyde	7.08	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

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280906.D Vial: 6
Acq On : 28 Jul 2009 9:54 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:29 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.69min 102.369ng/ml

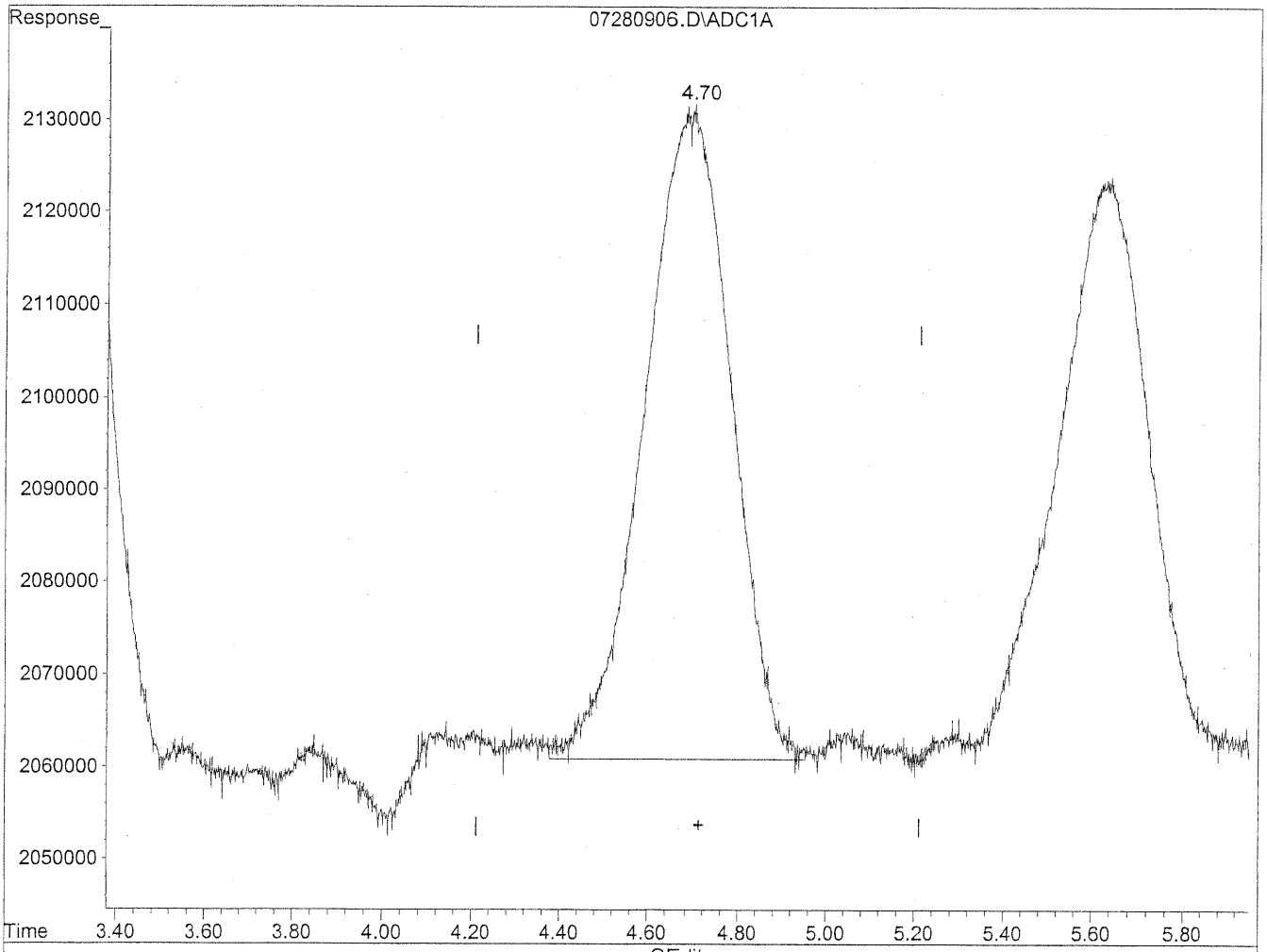
response 11317409

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280906.D Vial: 6
Acq On : 28 Jul 2009 9:54 am Operator: HC
Sample : 100ng/ml TO11A Std S21-07270905 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 15:29:52 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.70min 84.541ng/ml m

response 9346475

HC 7/28/09 1C

KG 7/29/09

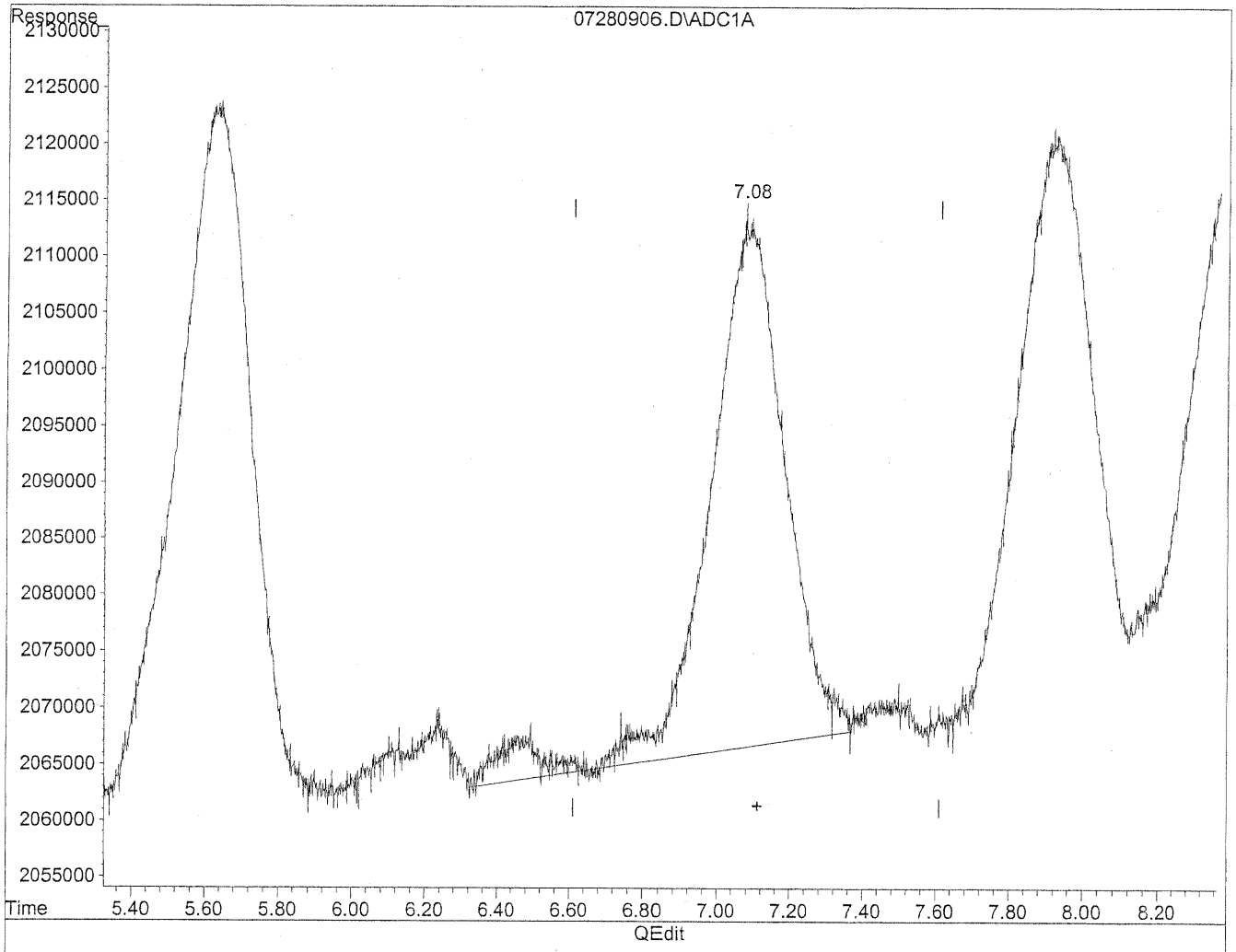
176

(+) = 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

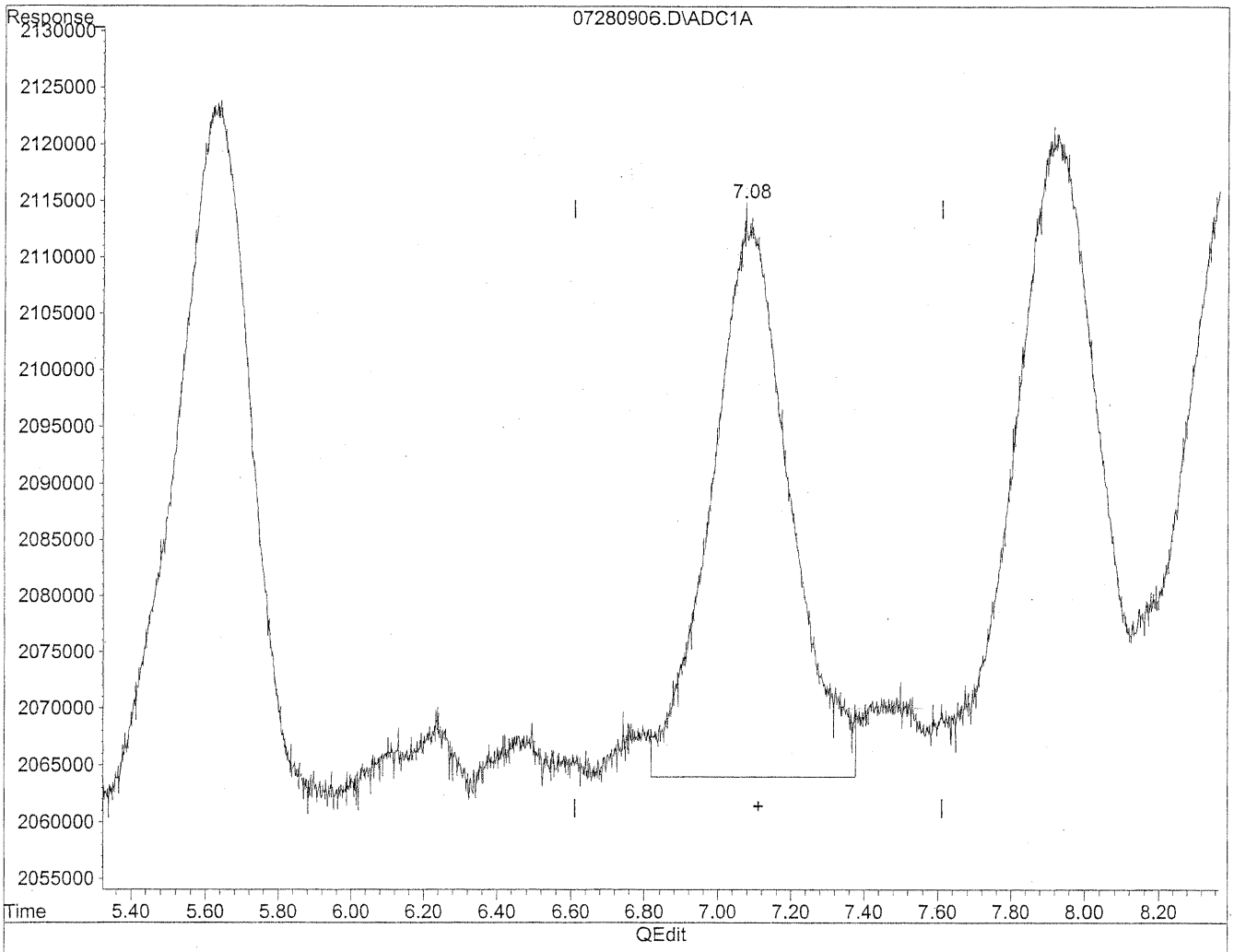
response 6819663

(+) = 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

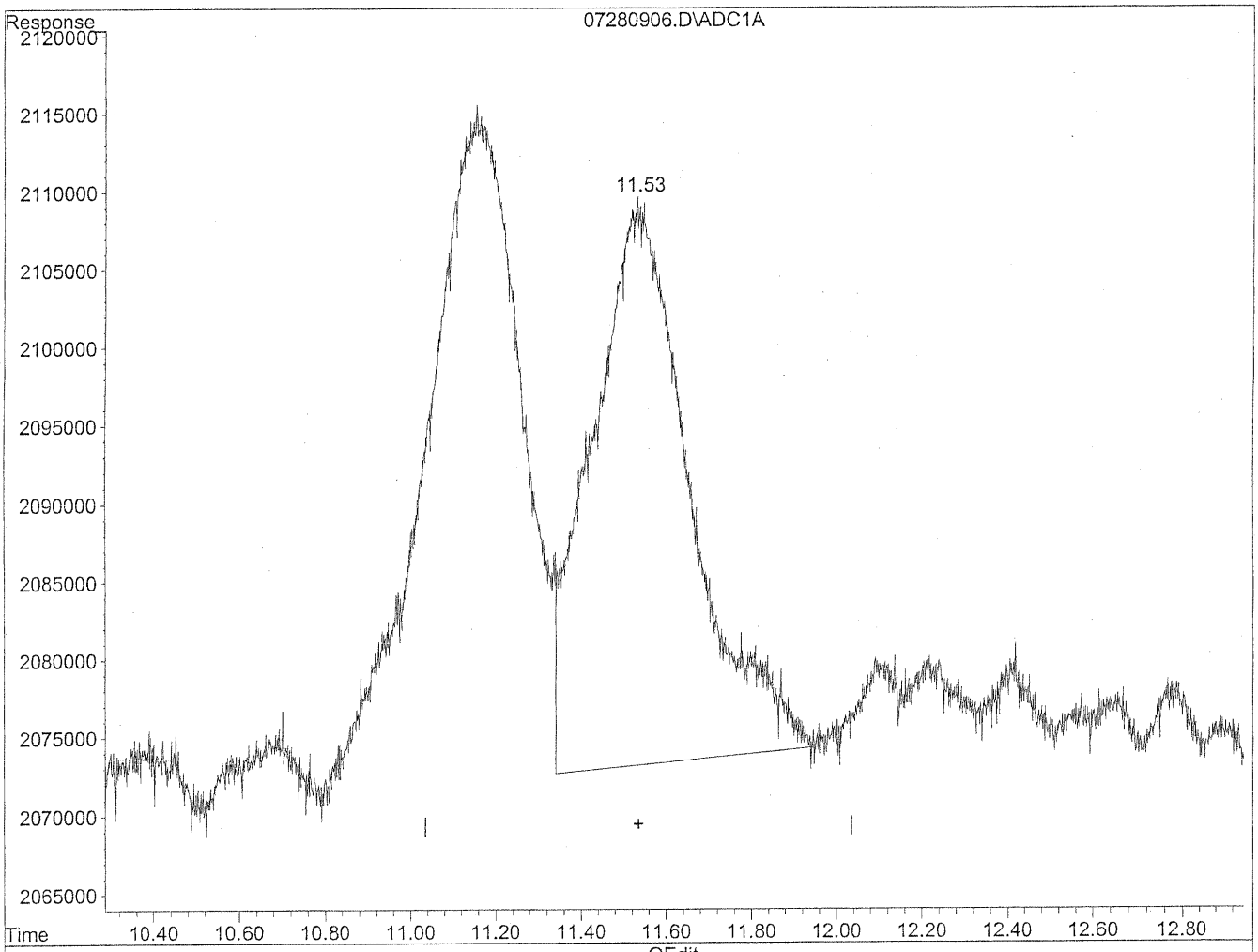
7/29/09

178

(+) = Expected Retention Time
07280906.D TO110709.M Tue Jul 28 14:30:30 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: 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

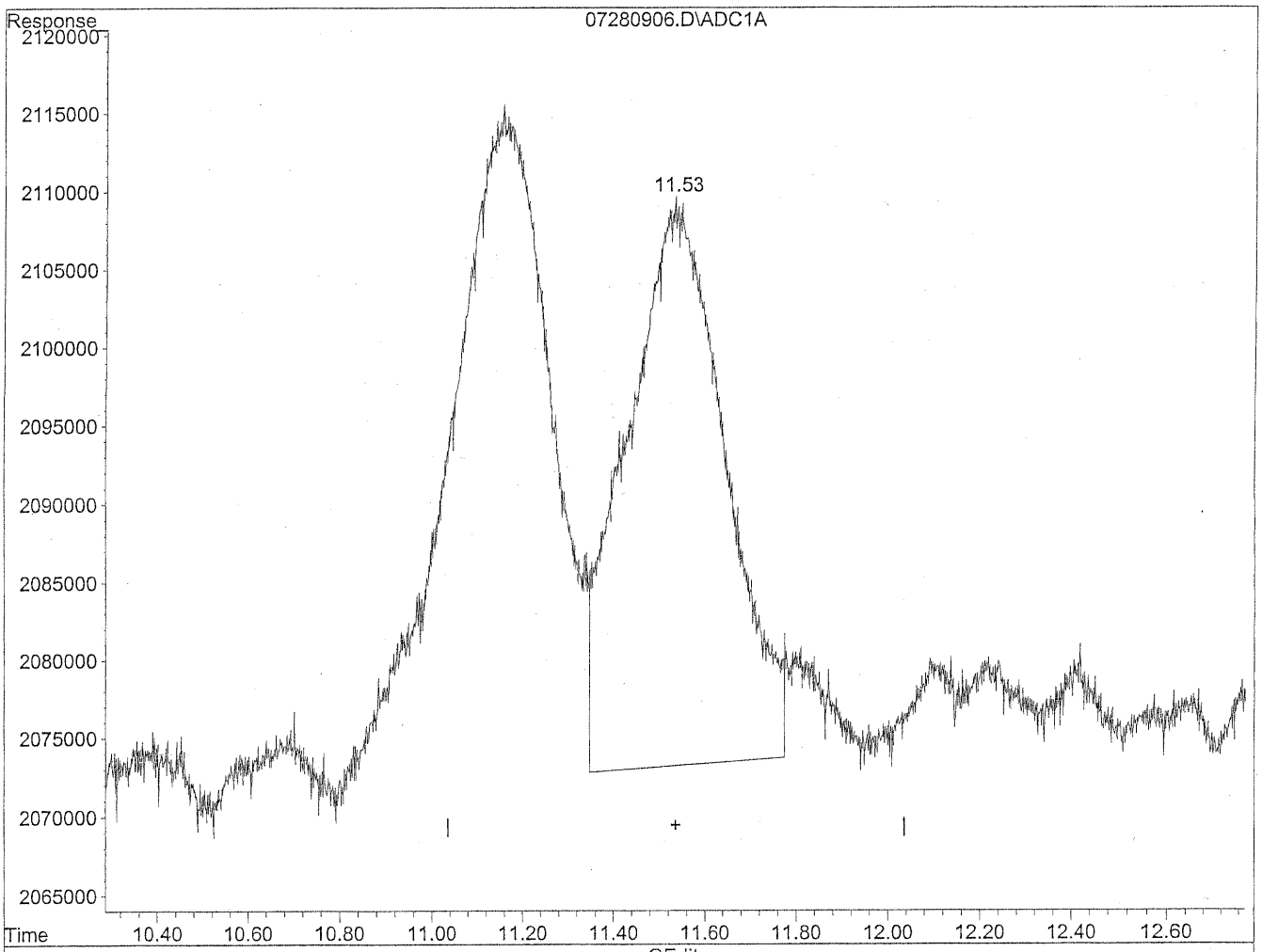
response 5798505

(+) = 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

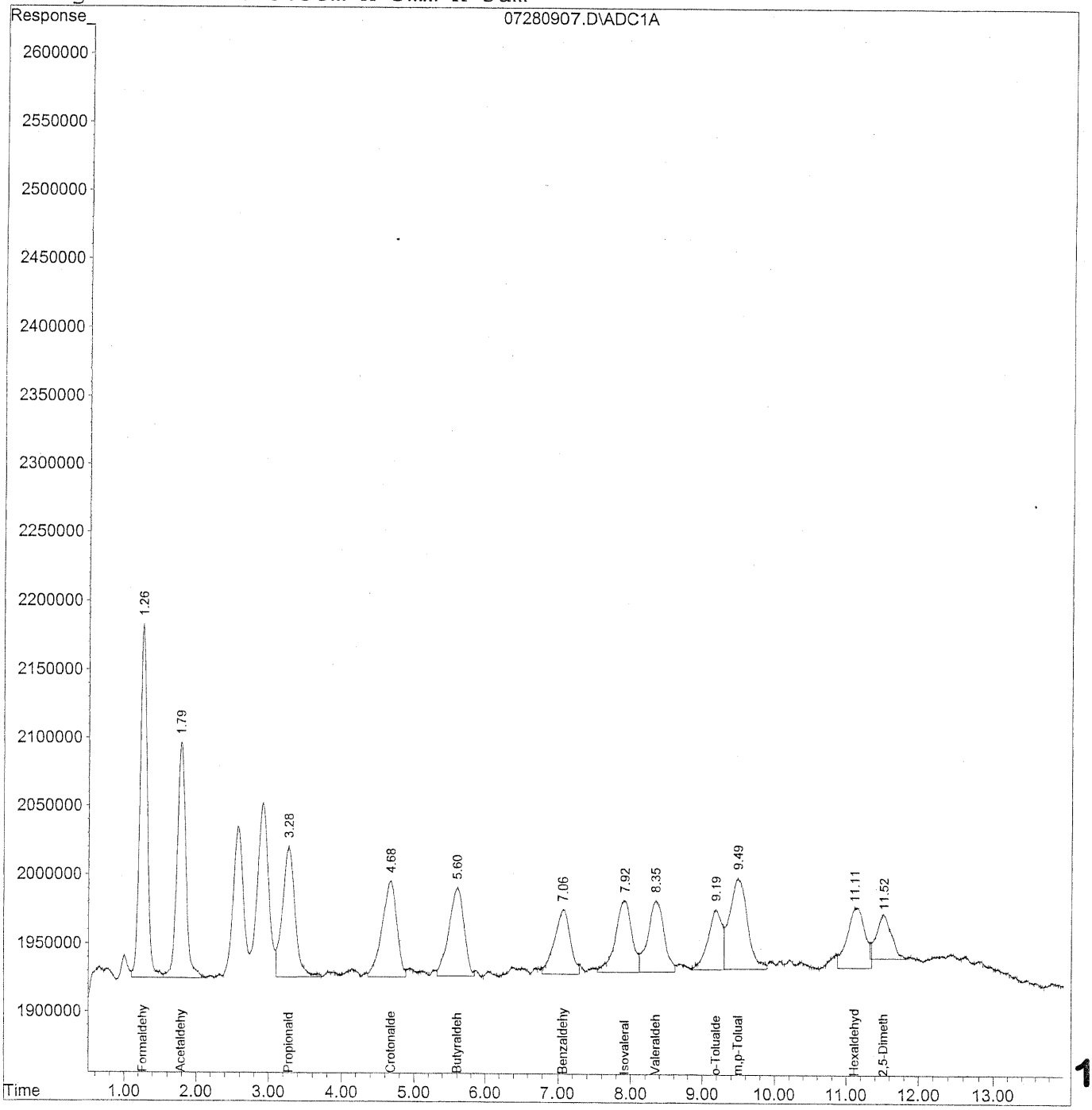
(+) = Expected Retention Time

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

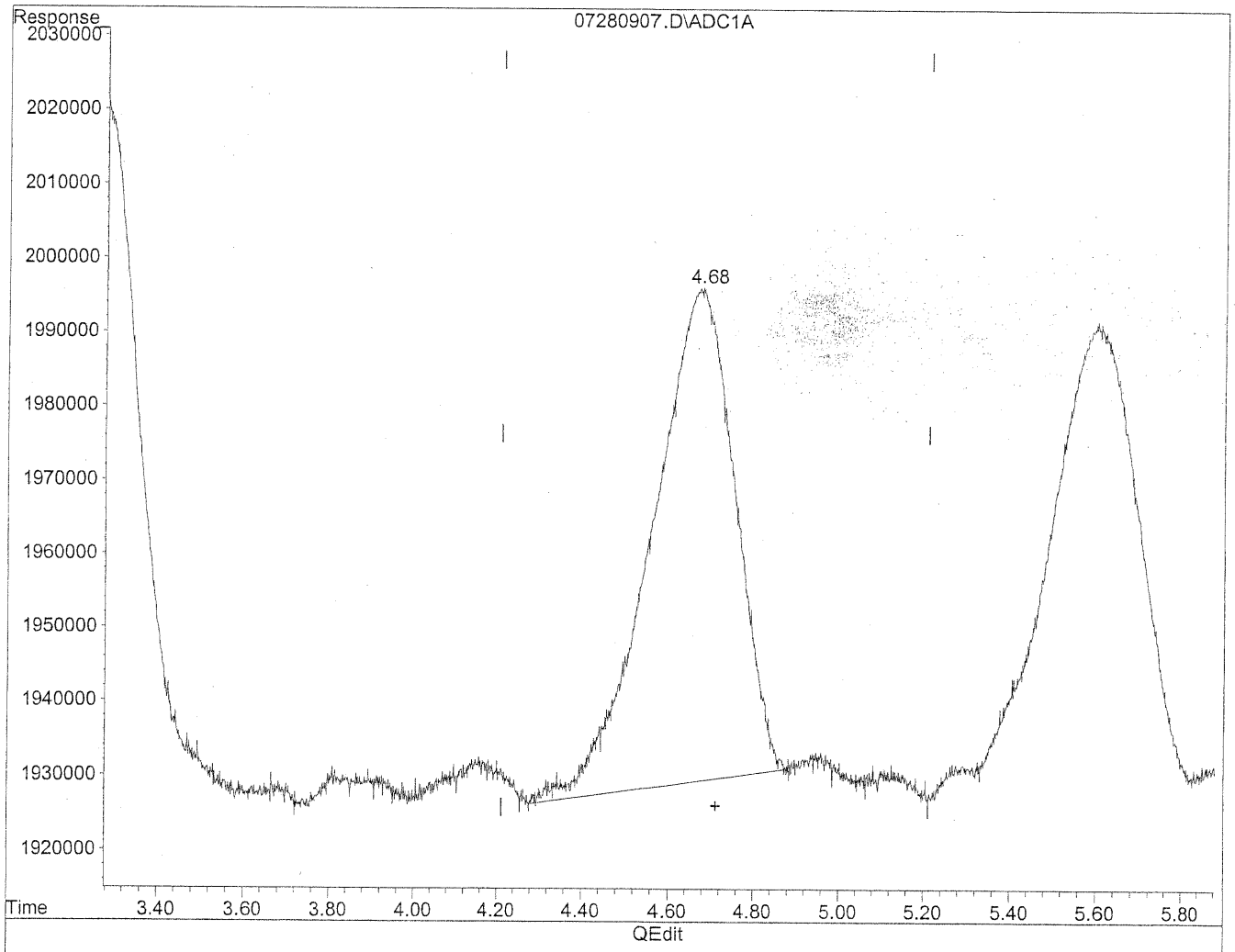
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.26	18449443	105.069	ng/ml
2)	Acetaldehyde	1.79	14434553	107.002	ng/ml
3)	Propionaldehyde	3.28	11389784	110.800	ng/ml
4)	Crotonaldehyde	4.68	9814490	88.774	ng/mlm
5)	Butyraldehyde	5.60	9432197	117.206	ng/mlm
6)	Benzaldehyde	7.06	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.
07280907.D TO110709.M Tue Jul 28 17:19:21 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.67min 80.883ng/ml

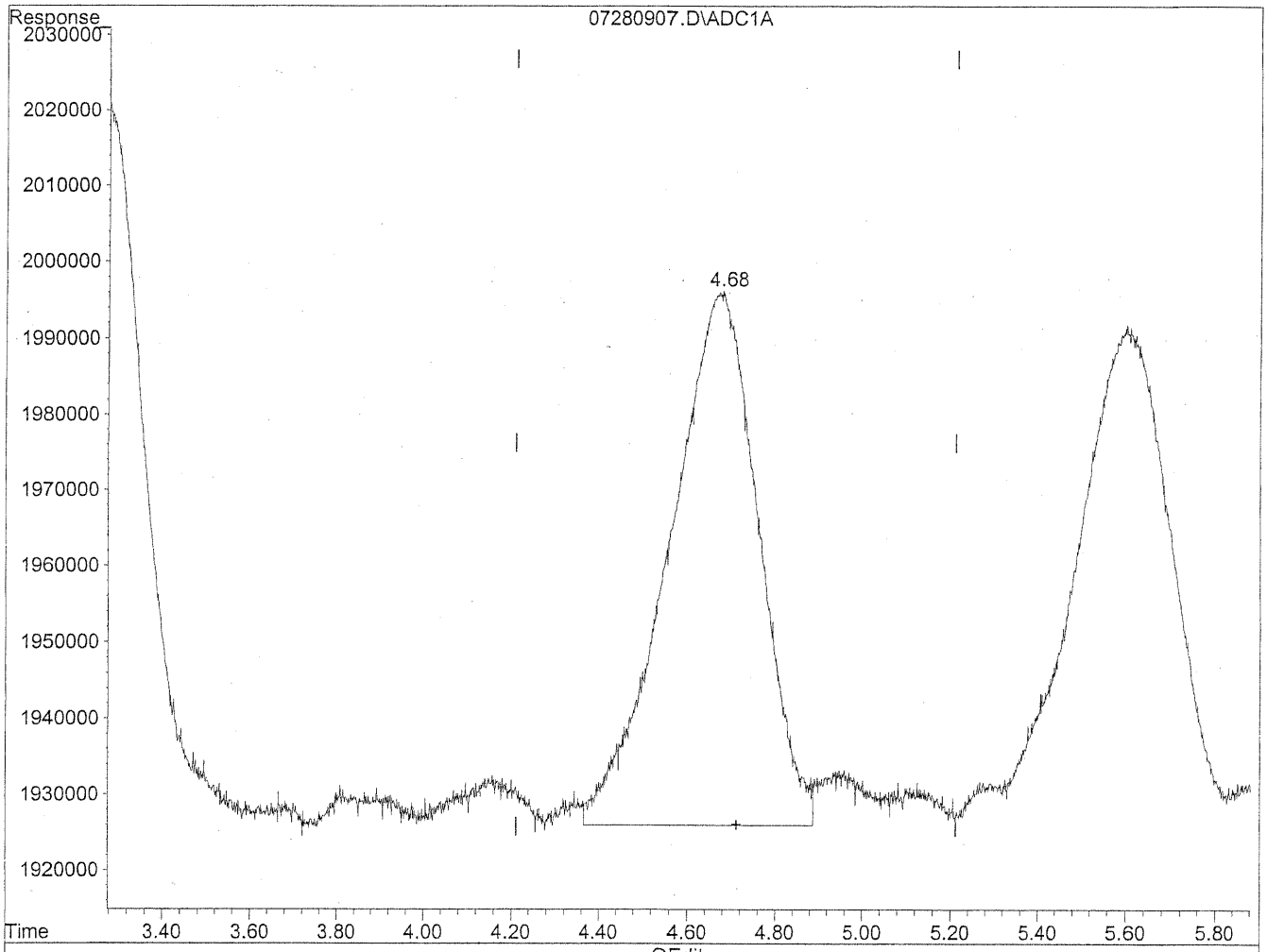
response 8942013

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.68min 88.774ng/ml m

response 9814490

AC 7/28/09 TC

KR 7/29/09

184

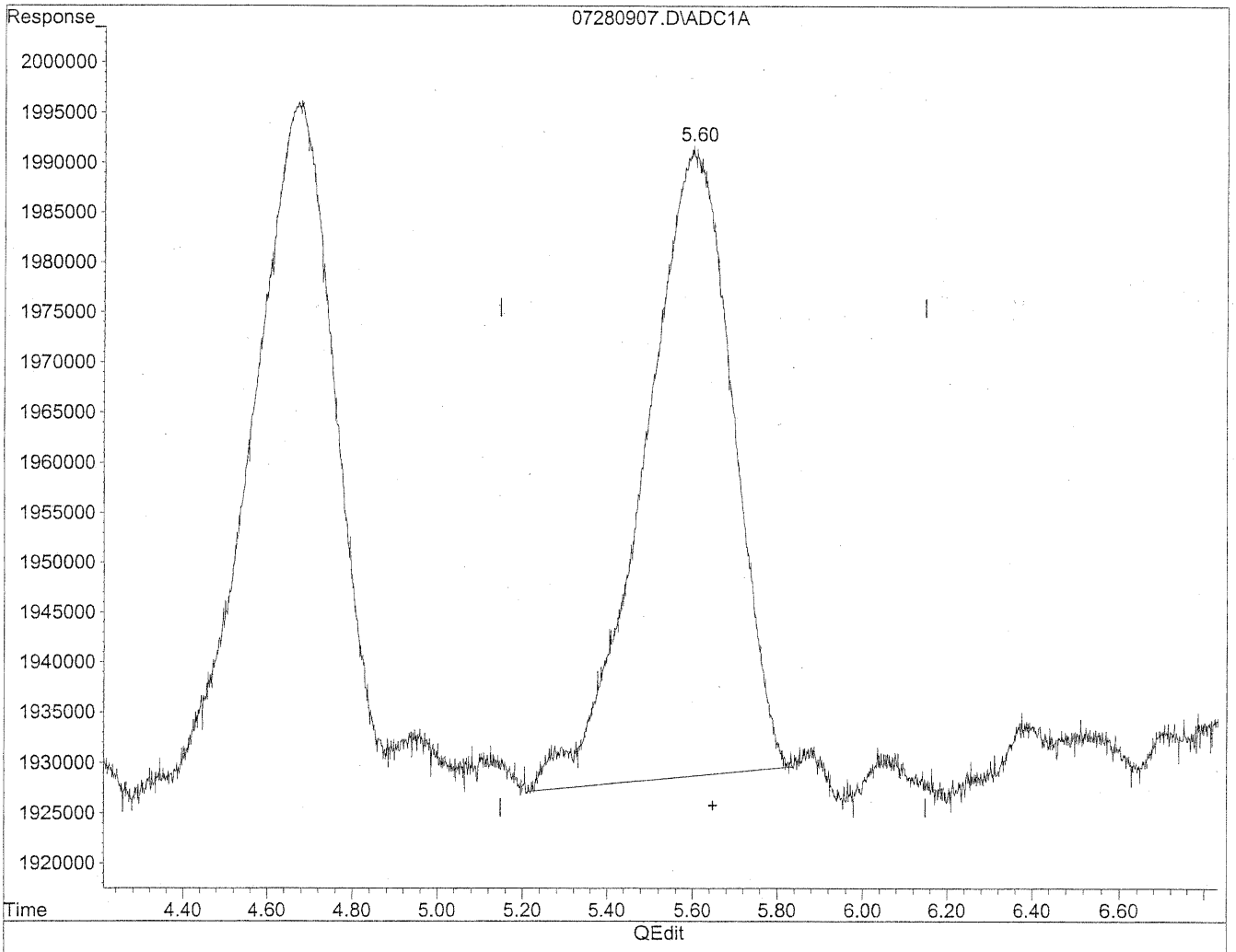
(+) = Expected Retention Time

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

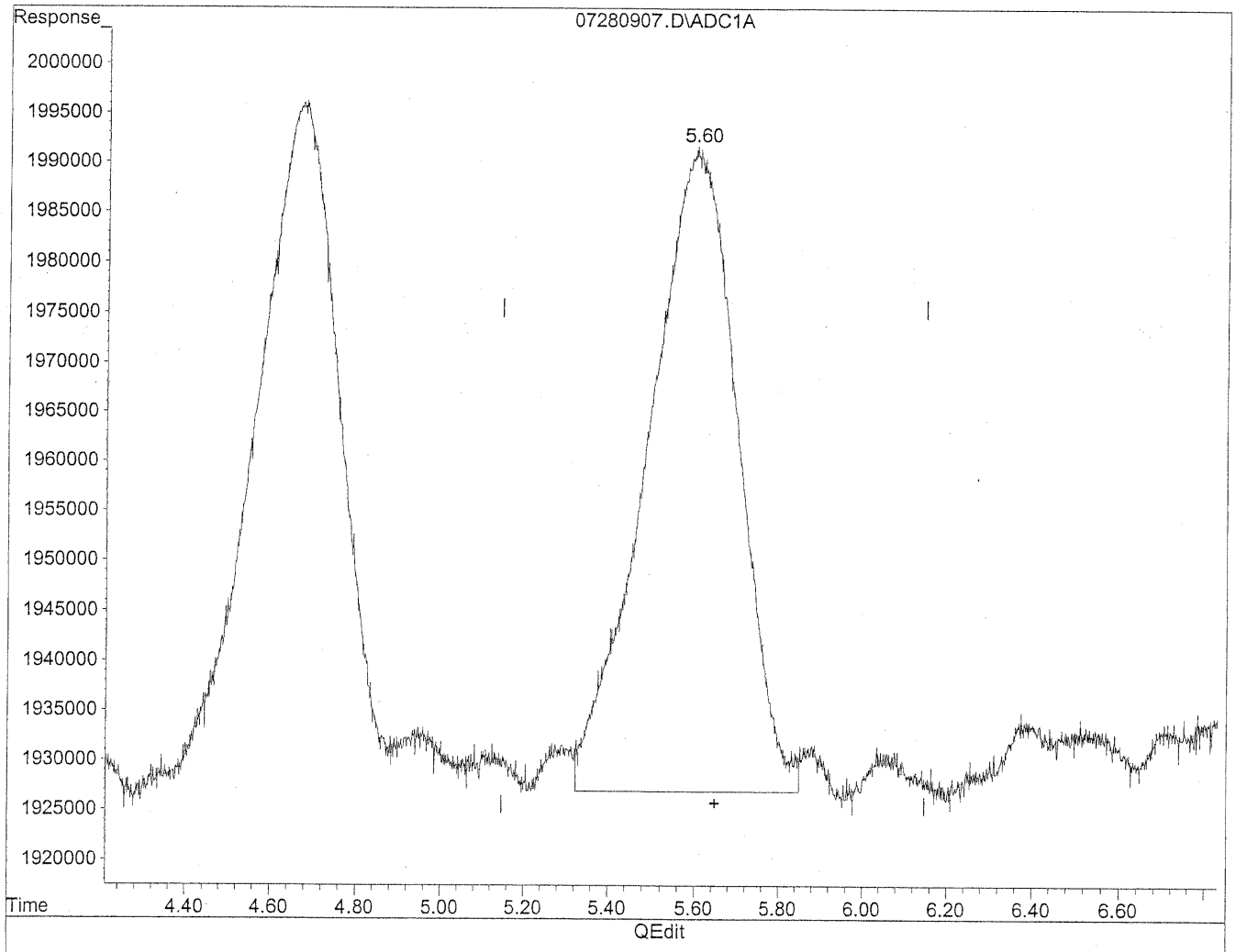
(+) = Expected Retention Time

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

HC 7/29/09

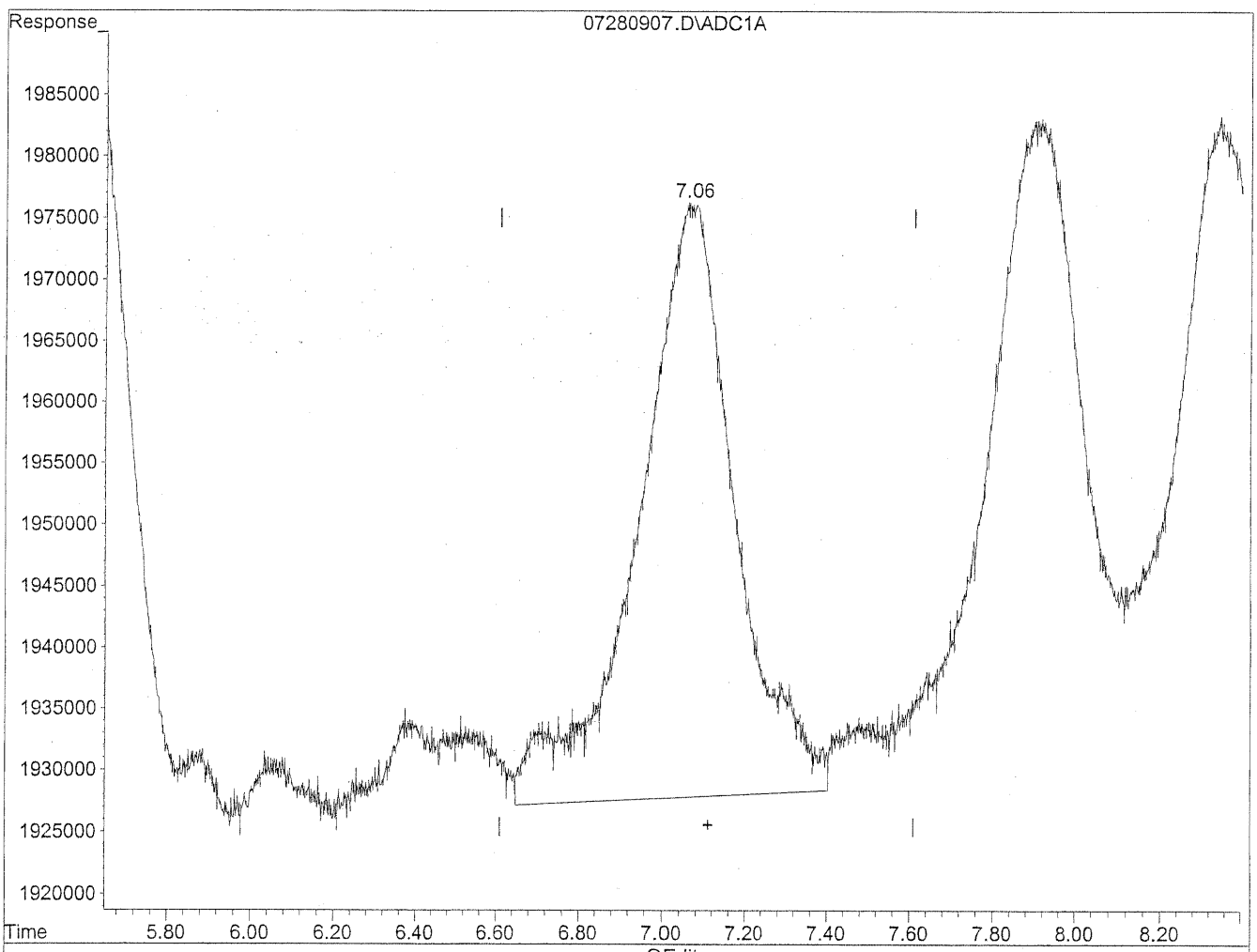
186

(+) = Expected Retention Time
07280907.D TO110709.M Tue Jul 28 14:31:25 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

7.07min 123.223ng/ml

response 7772036

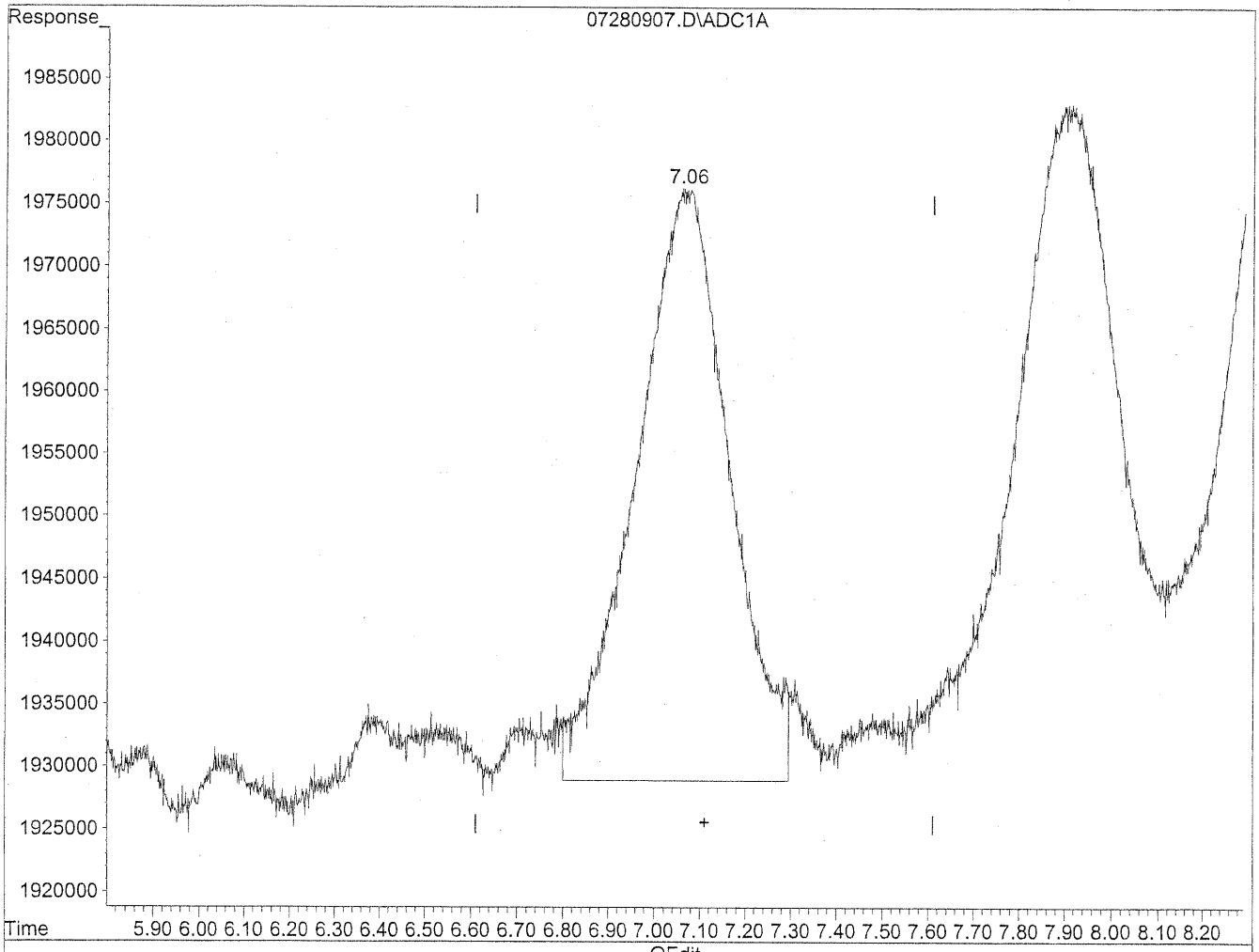
187

(+) = 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

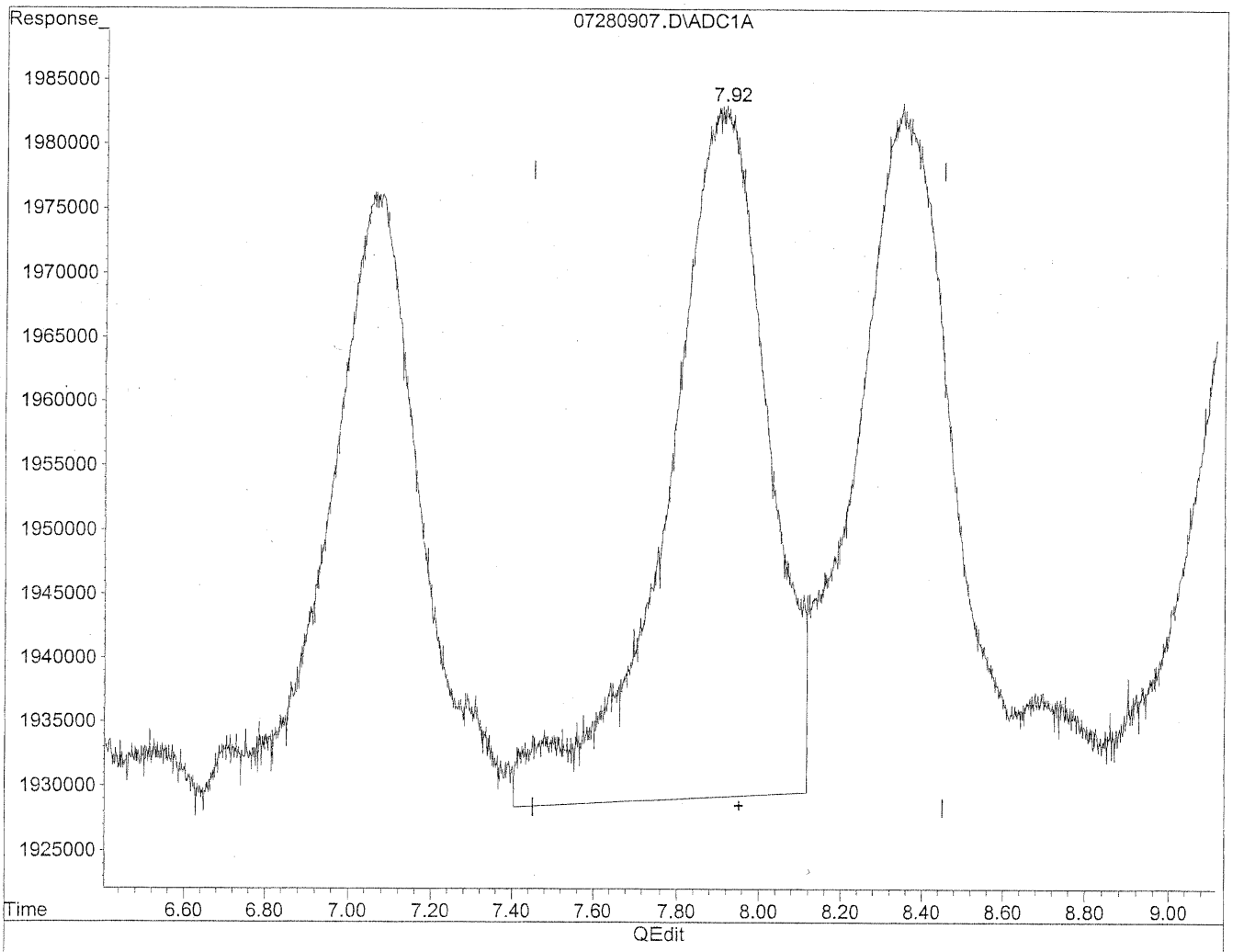
1427/24/09

(+) = 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

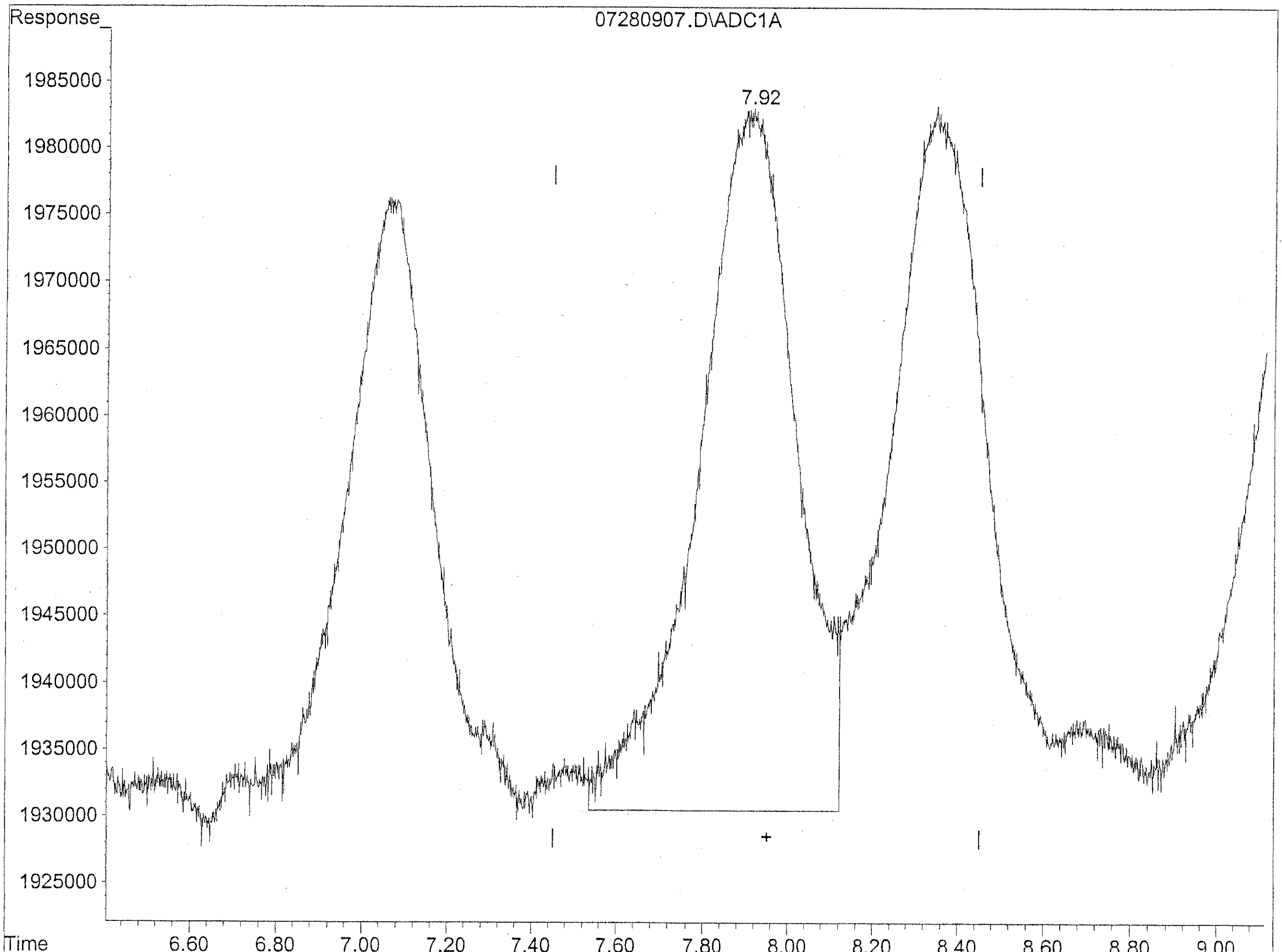
189

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



QEdit

(7) Isovaleraldehyde

7.92min 94.058ng/ml m

response 8338385

HC 7/28/09 IC

KR 7/29/09

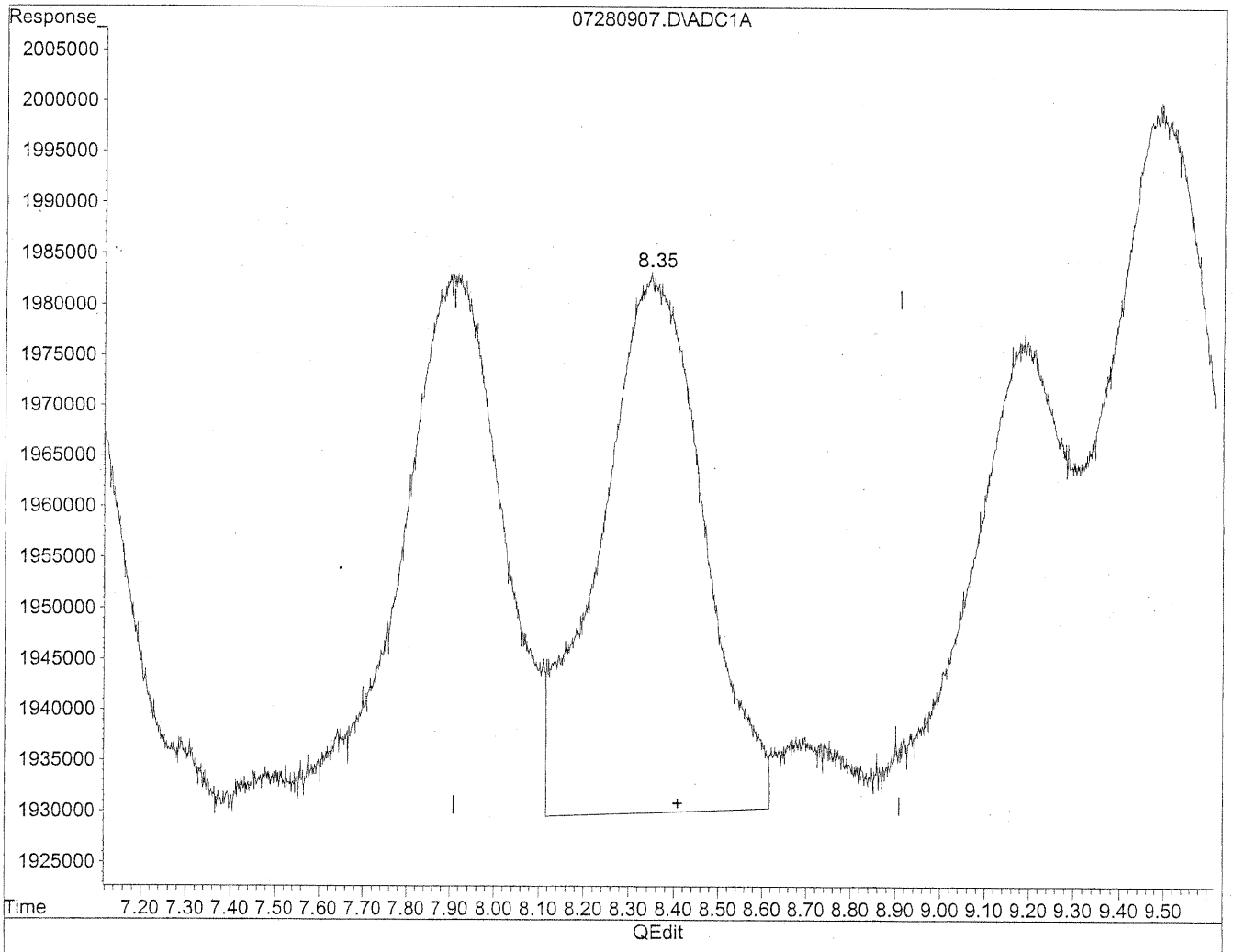
(+) = Expected Retention Time

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

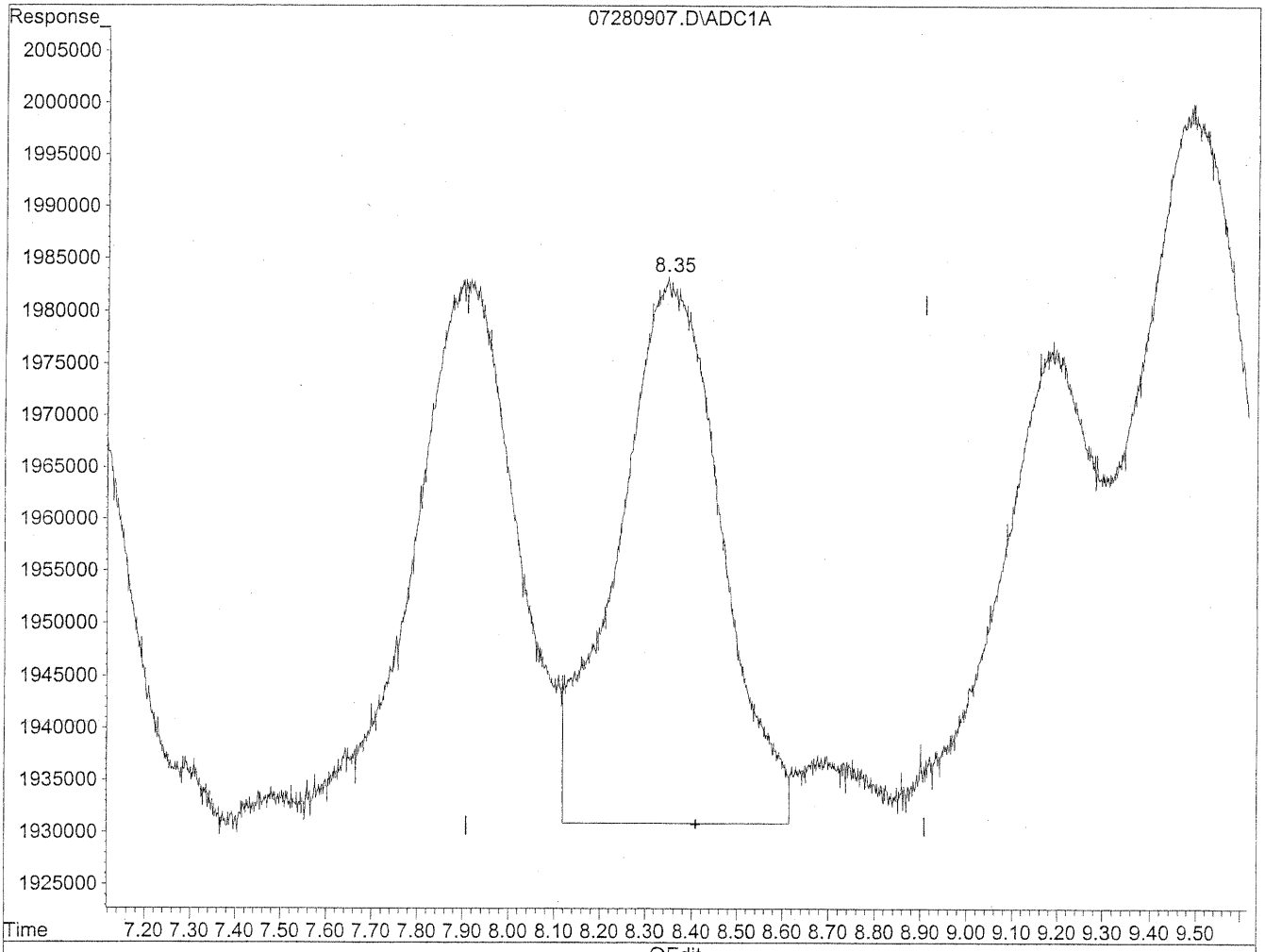
191

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



QEdit

(8) Valeraldehyde

8.35min 97.688ng/ml m

response 8117341

HC 7/28/09 RL

KF 7/29/09

192

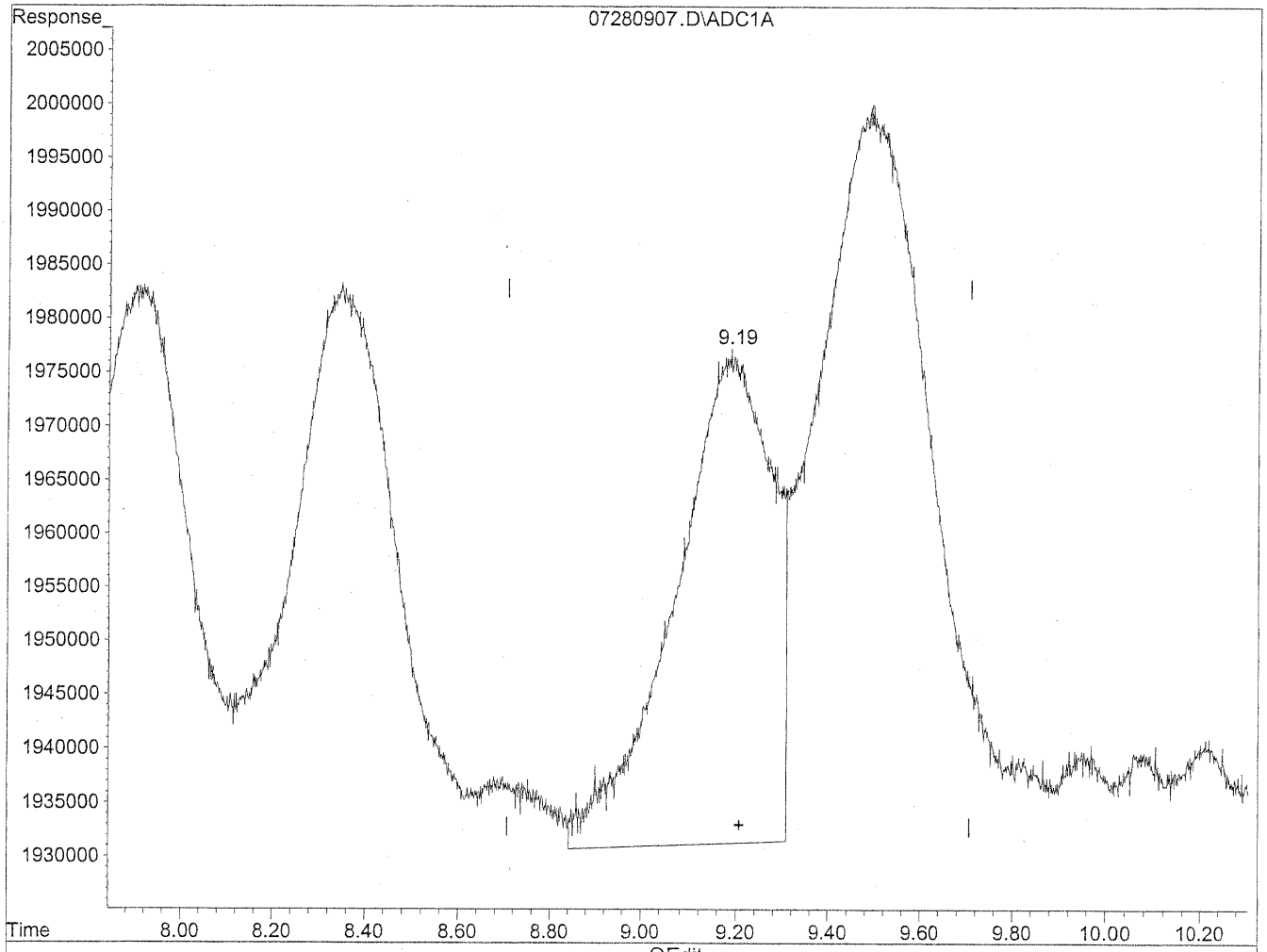
(+) = Expected Retention Time

07280907.D TO110709.M Tue Jul 28 14:32:16 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280907.D Vial: 7
Acq On : 28 Jul 2009 10:09 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:30 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

9.19min 121.312ng/ml

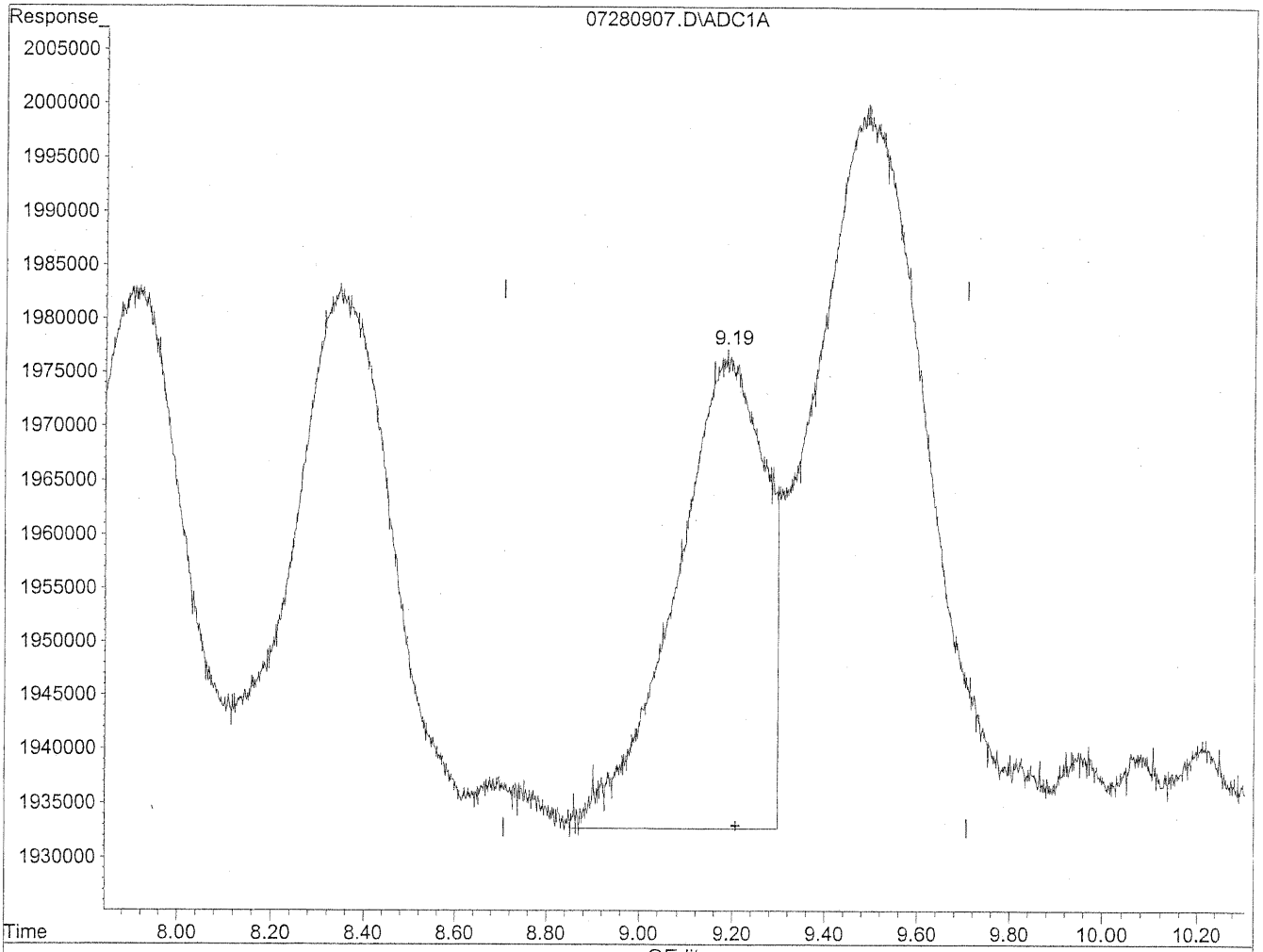
response 6535124

(+) = Expected Retention Time

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

7/28/09

KR 7/29/09

194

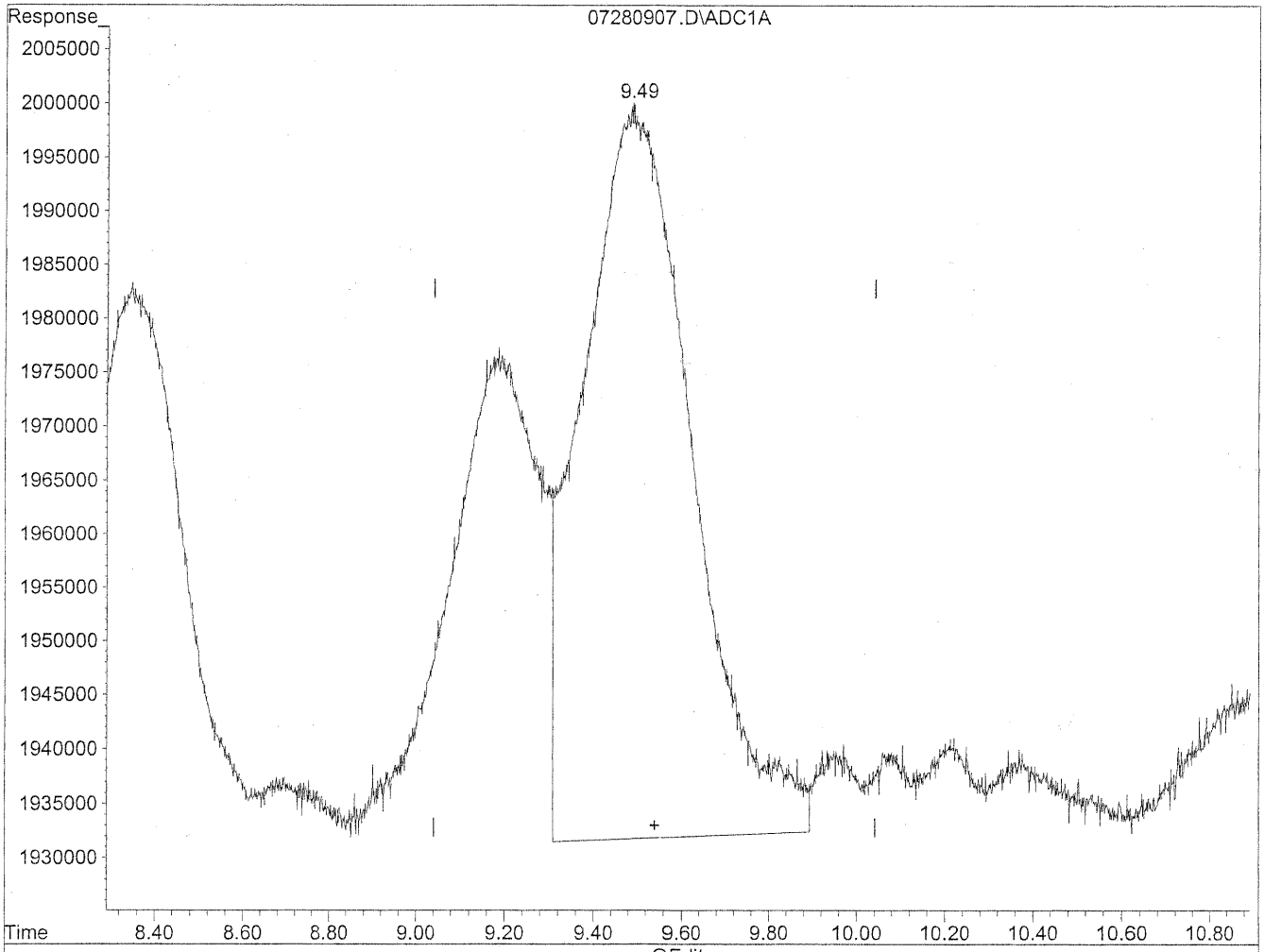
(+) = Expected Retention Time

07280907.D TO110709.M Tue Jul 28 14:32:36 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



(10) m,p-Tolualdehyde

9.49min 217.917ng/ml

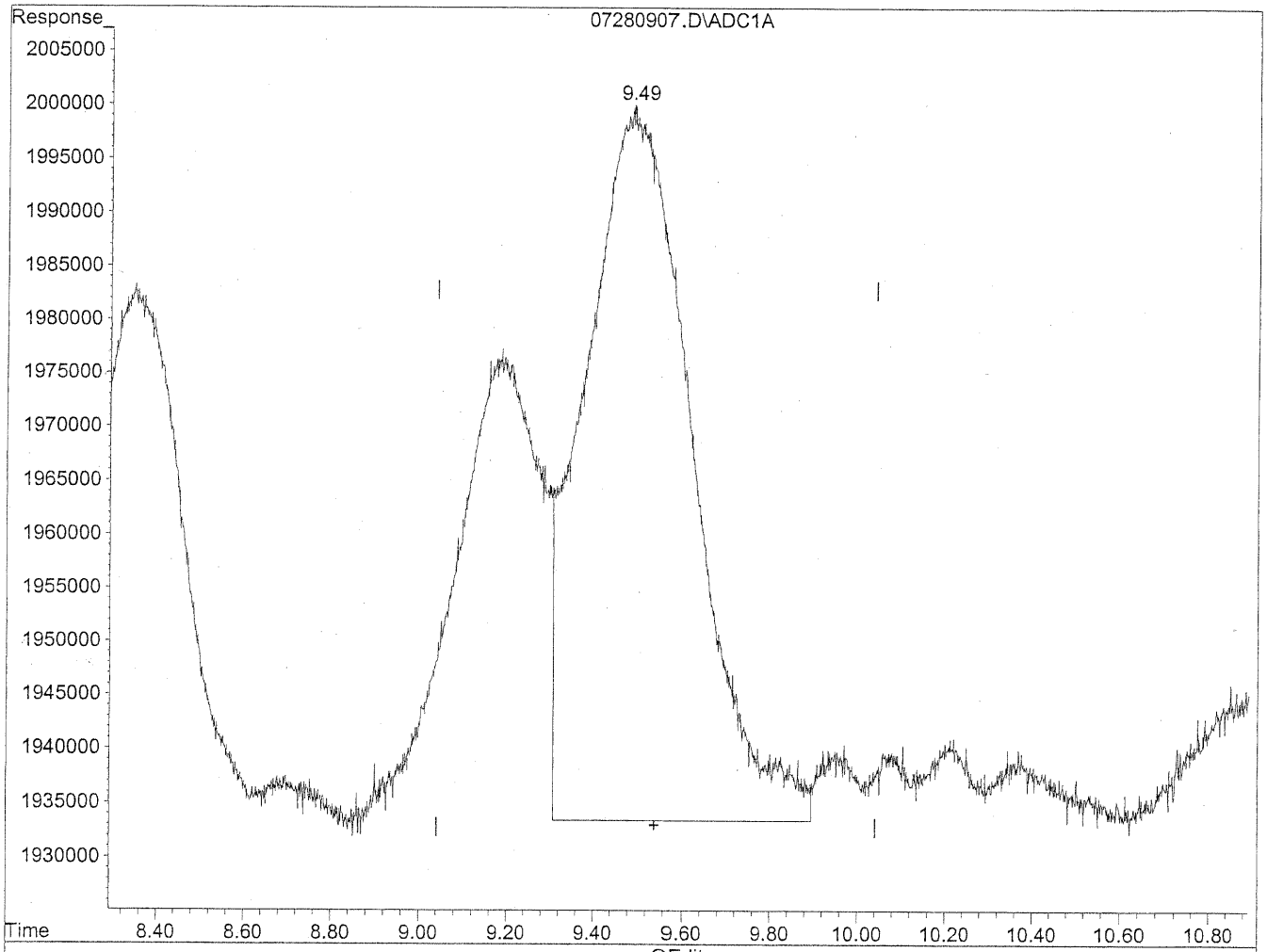
response 11738041

(+) = 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



QEdit

(10) m,p-Tolualdehyde

9.49min 208.581ng/ml m

response 11235135

HC 7/28/09

KR 7/29/09

196

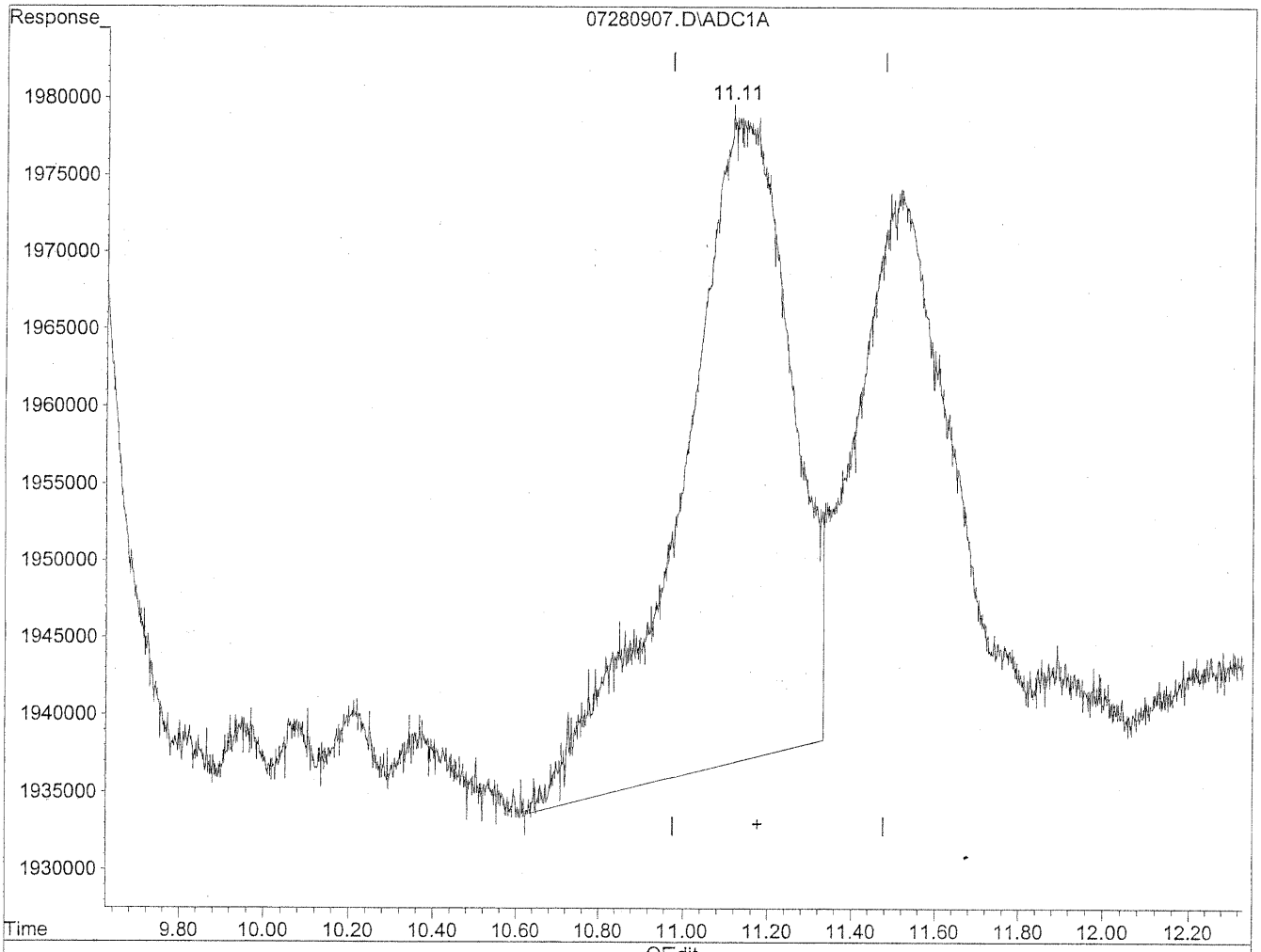
(+) = Expected Retention Time

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

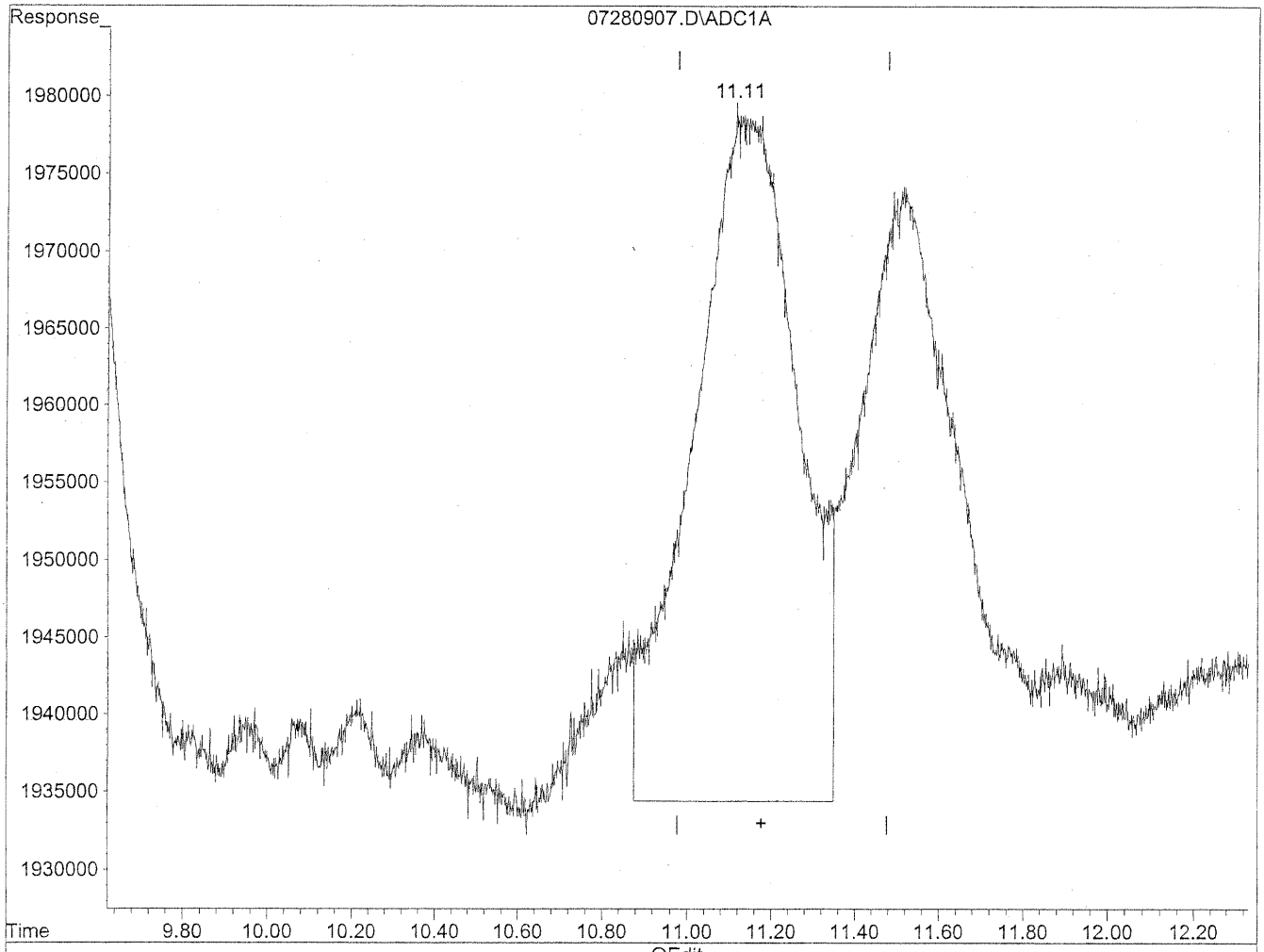
response 7552544

(+) = 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

(+) = Expected Retention Time

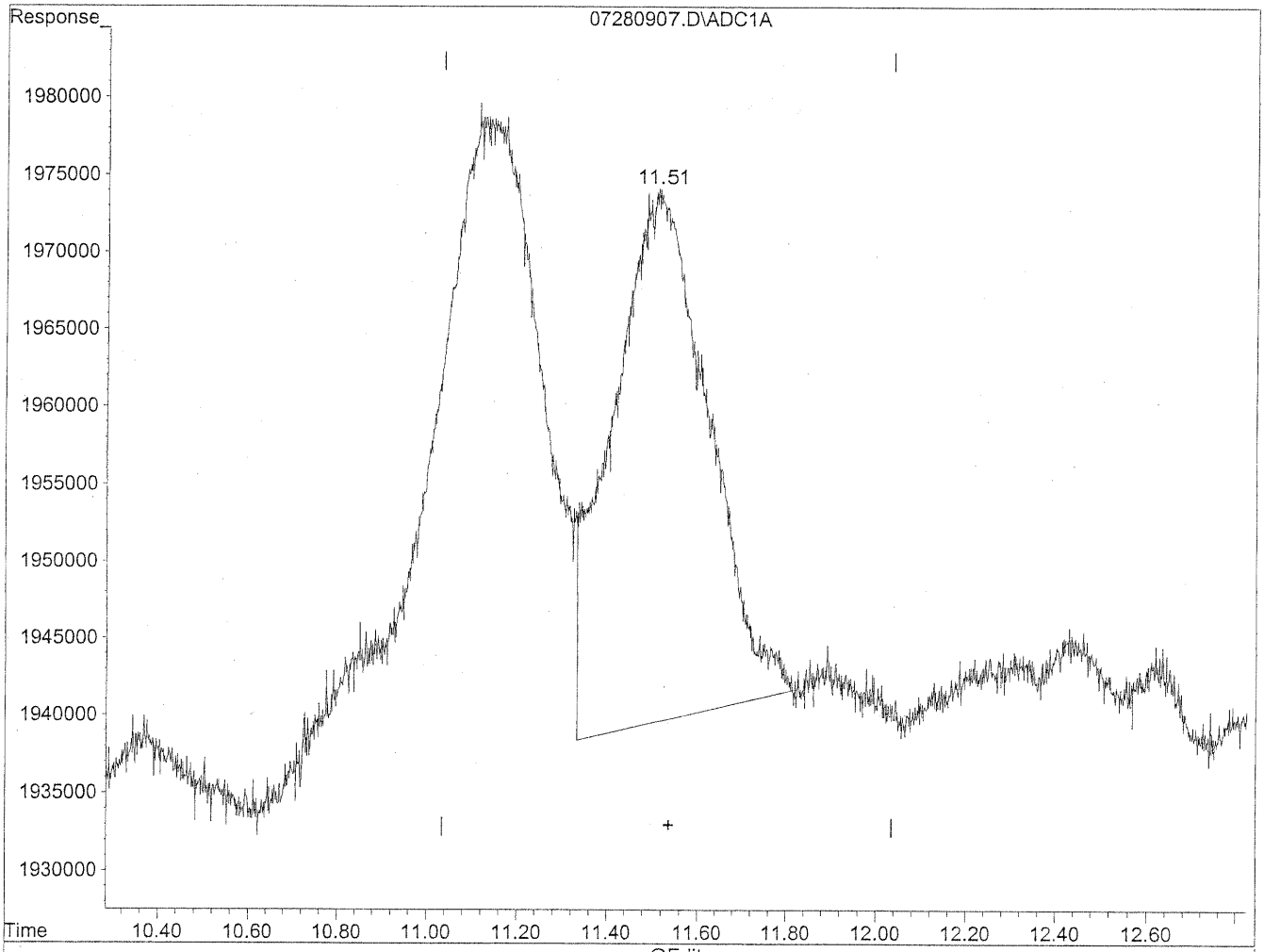
07280907.D TO110709.M Tue Jul 28 14:33:16 2009

198

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

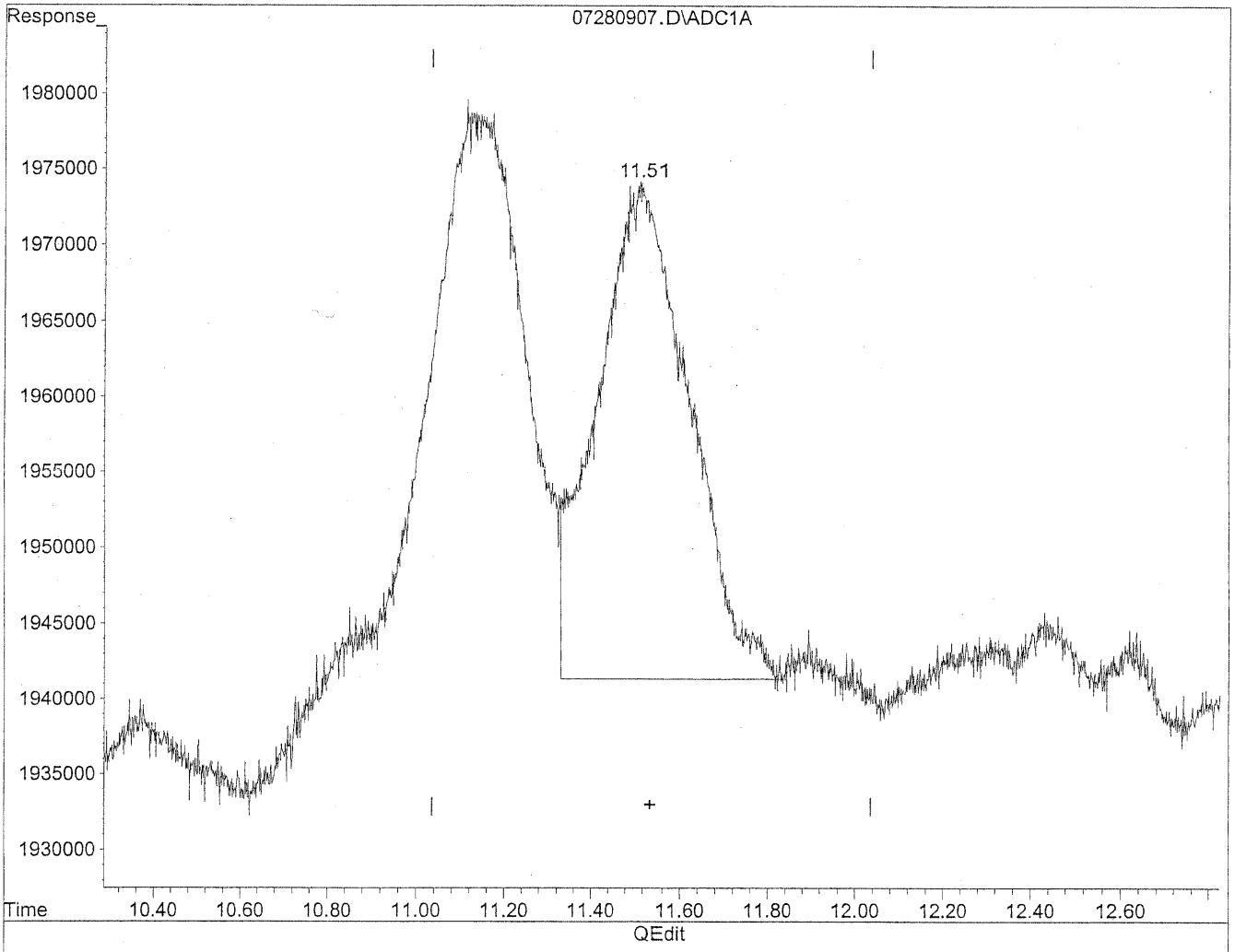
response 5084888

(+) = 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 PL

KR 7/29/09

200

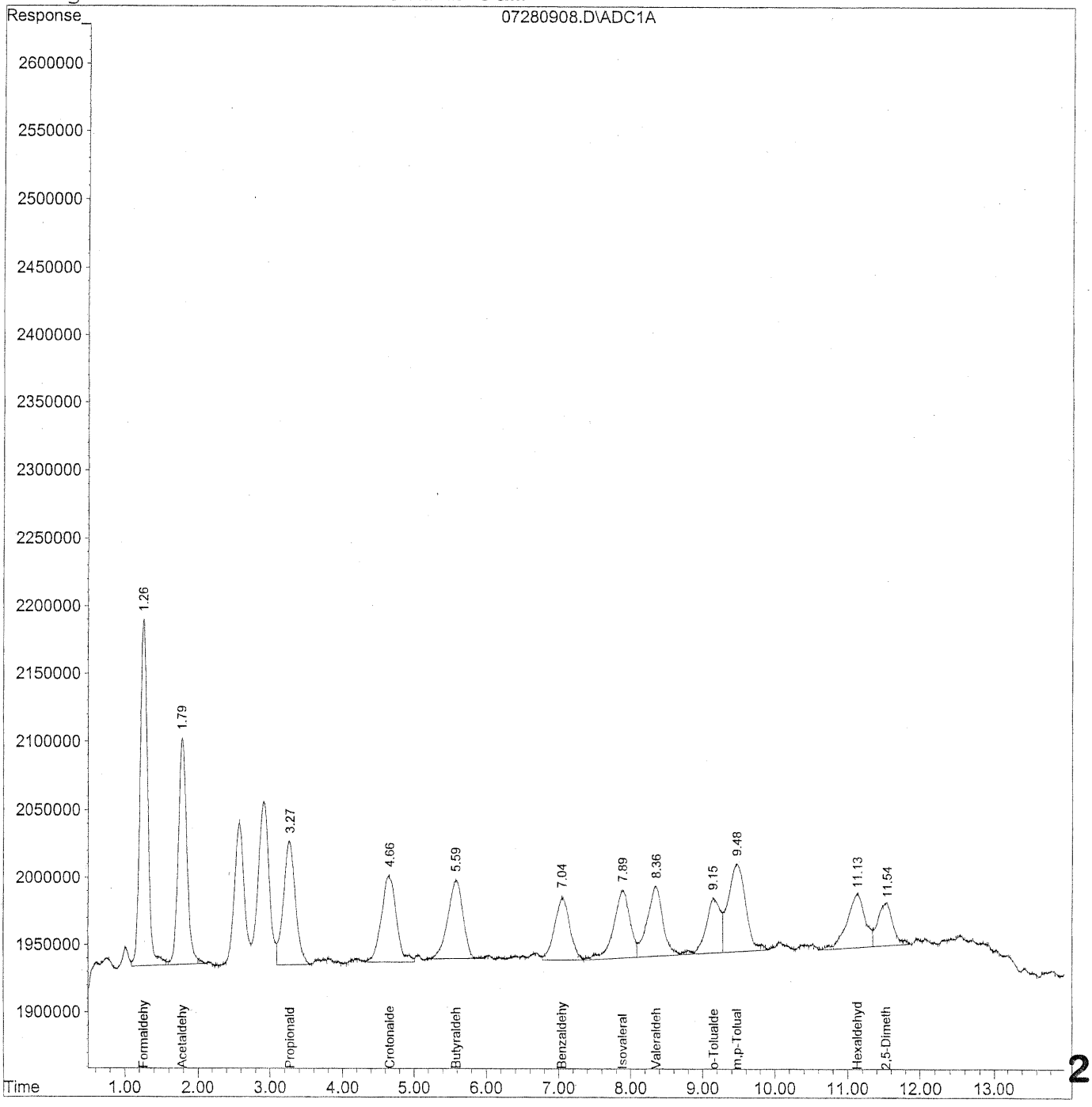
(+) = Expected Retention Time

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



201

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

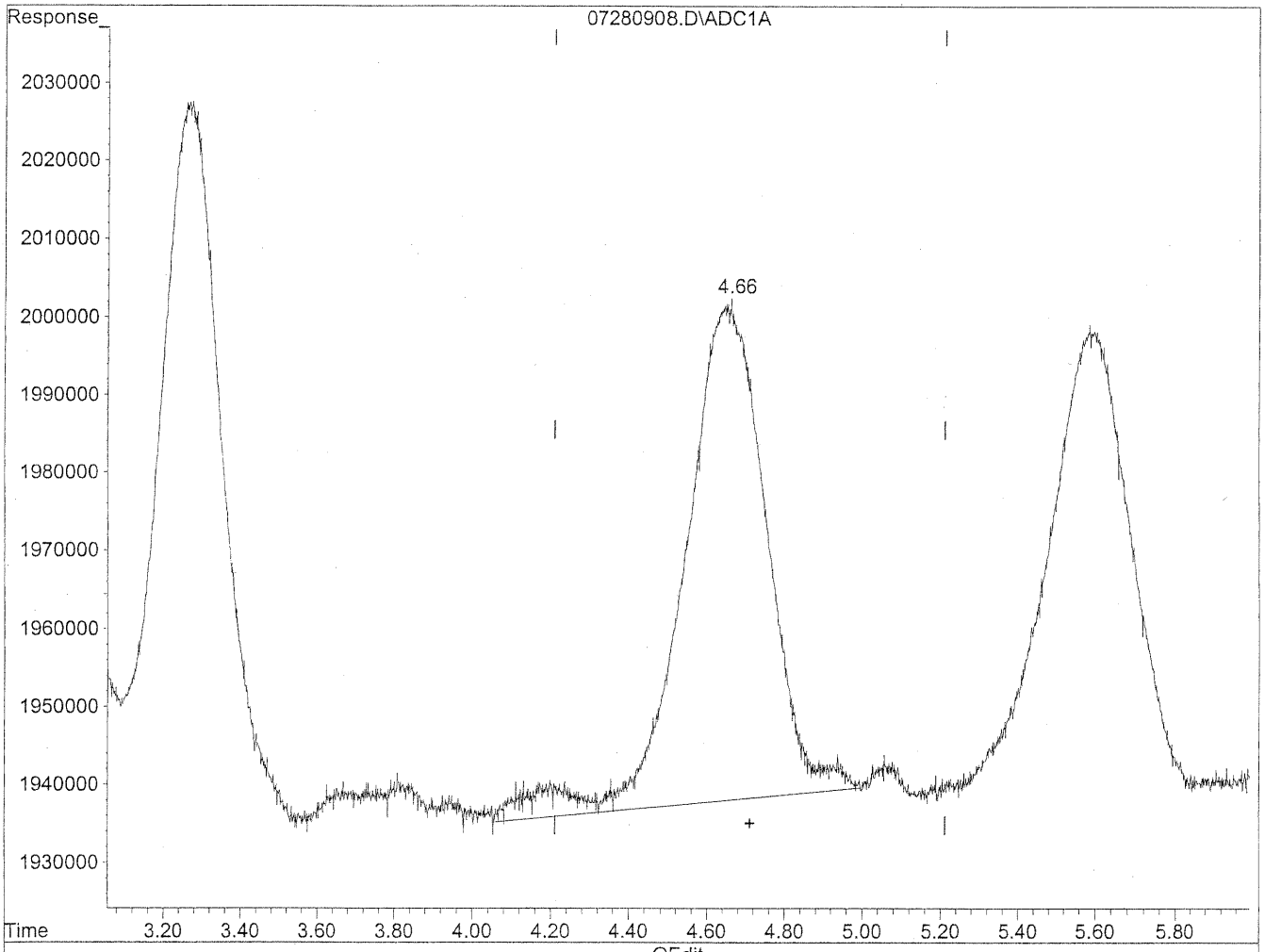
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.26	18400032	104.788	ng/ml
2)	Acetaldehyde	1.79	13737532	101.835	ng/ml
3)	Propionaldehyde	3.27	10633406	103.442	ng/ml
4)	Crotonaldehyde	4.66	9424529	85.247	ng/mlm
5)	Butyraldehyde	5.59	8463028	105.163	ng/ml
6)	Benzaldehyde	7.04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

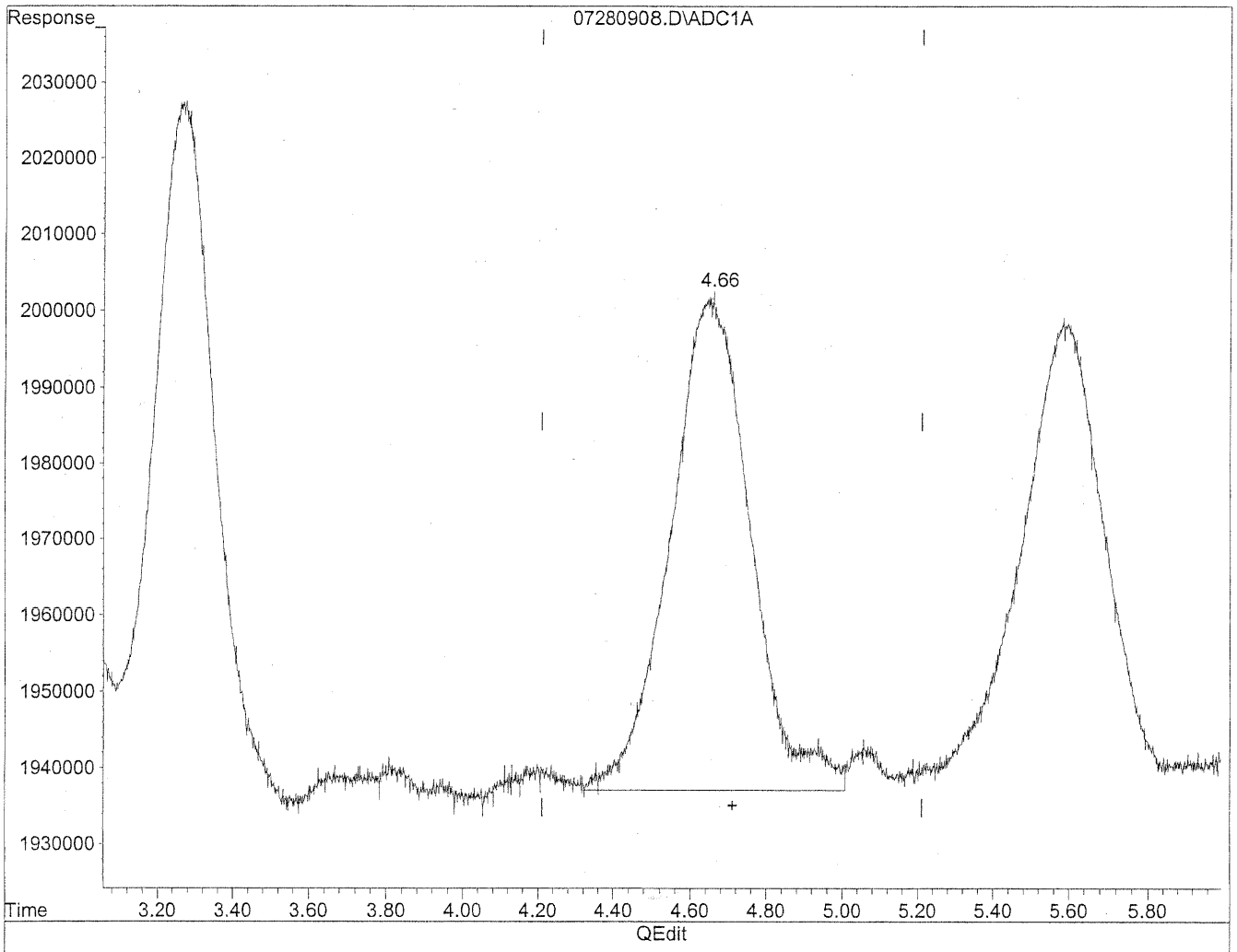
(+) = Expected Retention Time

07280908.D TO110709.M Tue Jul 28 14:33:54 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 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 STH

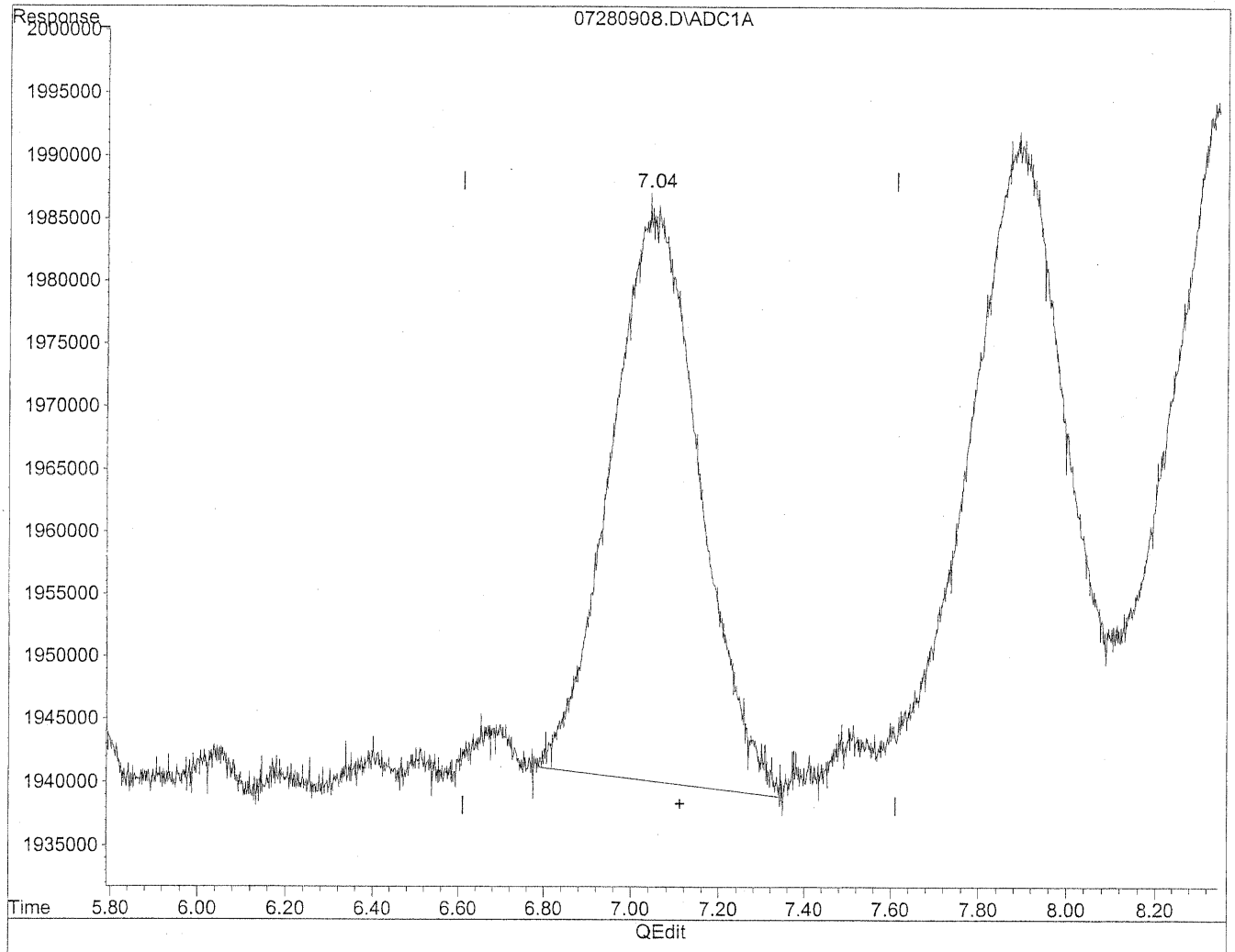
KF 7/29/09

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280908.D Vial: 8
Acq On : 28 Jul 2009 10:24 am Operator: HC
Sample : 100ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:33 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(6) Benzaldehyde

7.05min 101.515ng/ml

response 6402857

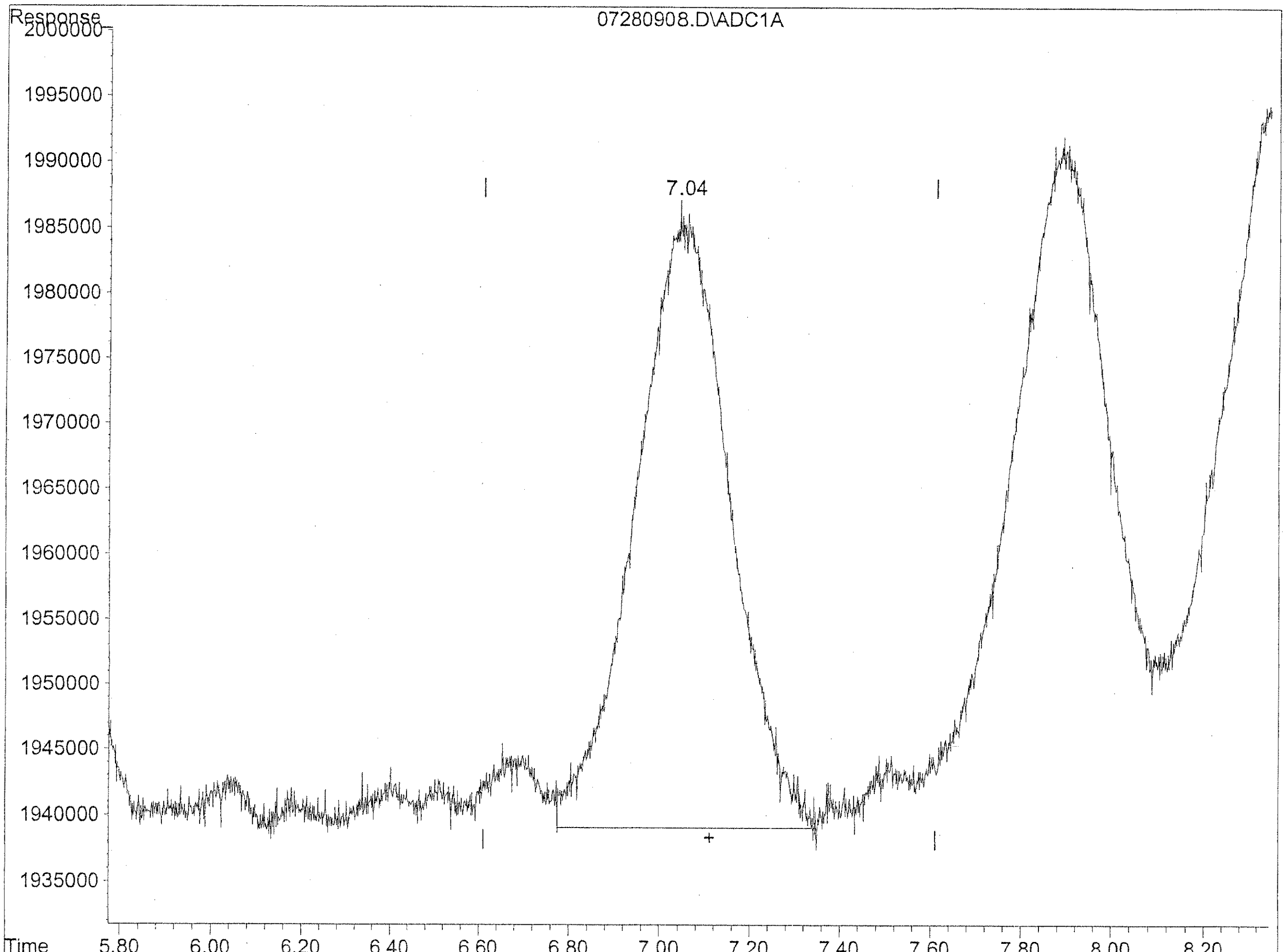
(+) = Expected Retention Time

07280908.D TO110709.M Tue Jul 28 14:34:27 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280908.D Vial: 8
Acq On : 28 Jul 2009 10:24 am Operator: HC
Sample : 100ng/ml TO11A Std S21-07270905 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:34 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 15:29:52 2009
Response via : Multiple Level Calibration



QEdit

(6) Benzaldehyde

7.04min 106.795ng/ml m

response 6735919

HC 7/28/09 BC

KR 7/29/09

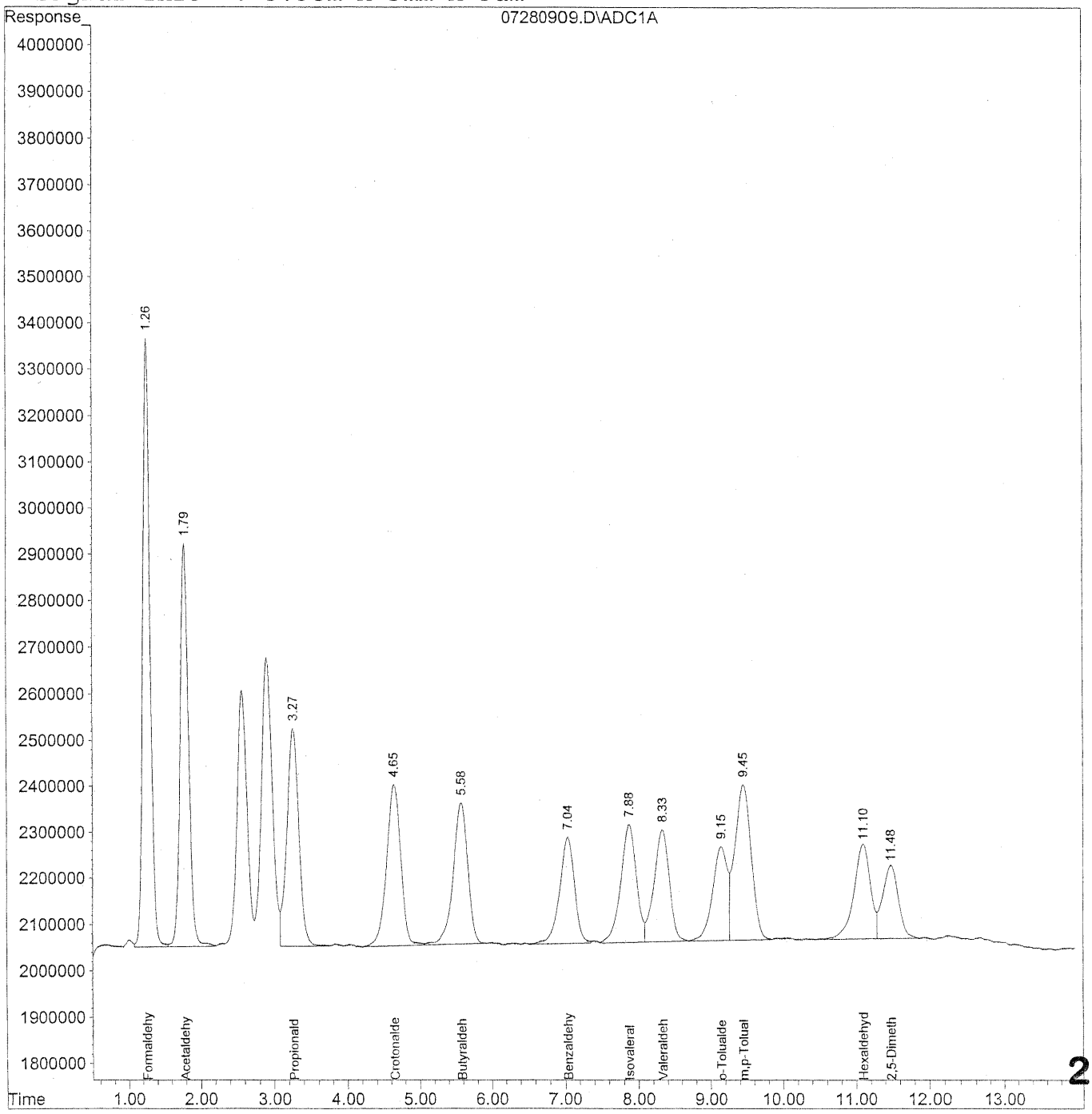
(+) = Expected Retention Time
07280908.D TO110709.M Tue Jul 28 17:38:30 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280909.D Vial: 9
Acq On : 28 Jul 2009 10:39 am Operator: HC
Sample : 500ng/ml TO11A Std S21-07270904 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:40 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

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



207

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

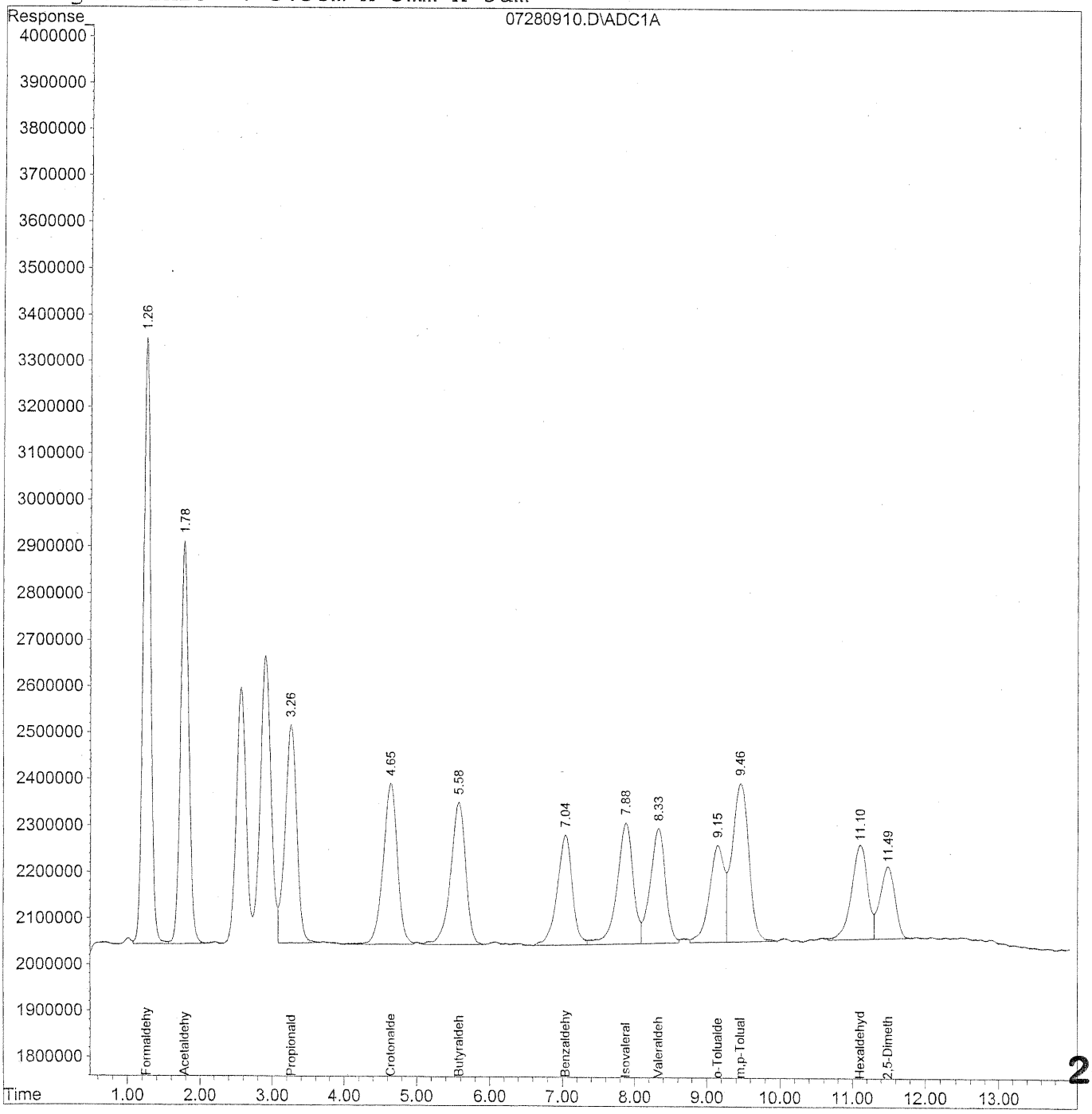
(f)=RT Delta > 1/2 Window (m)=manual int.

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



209

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

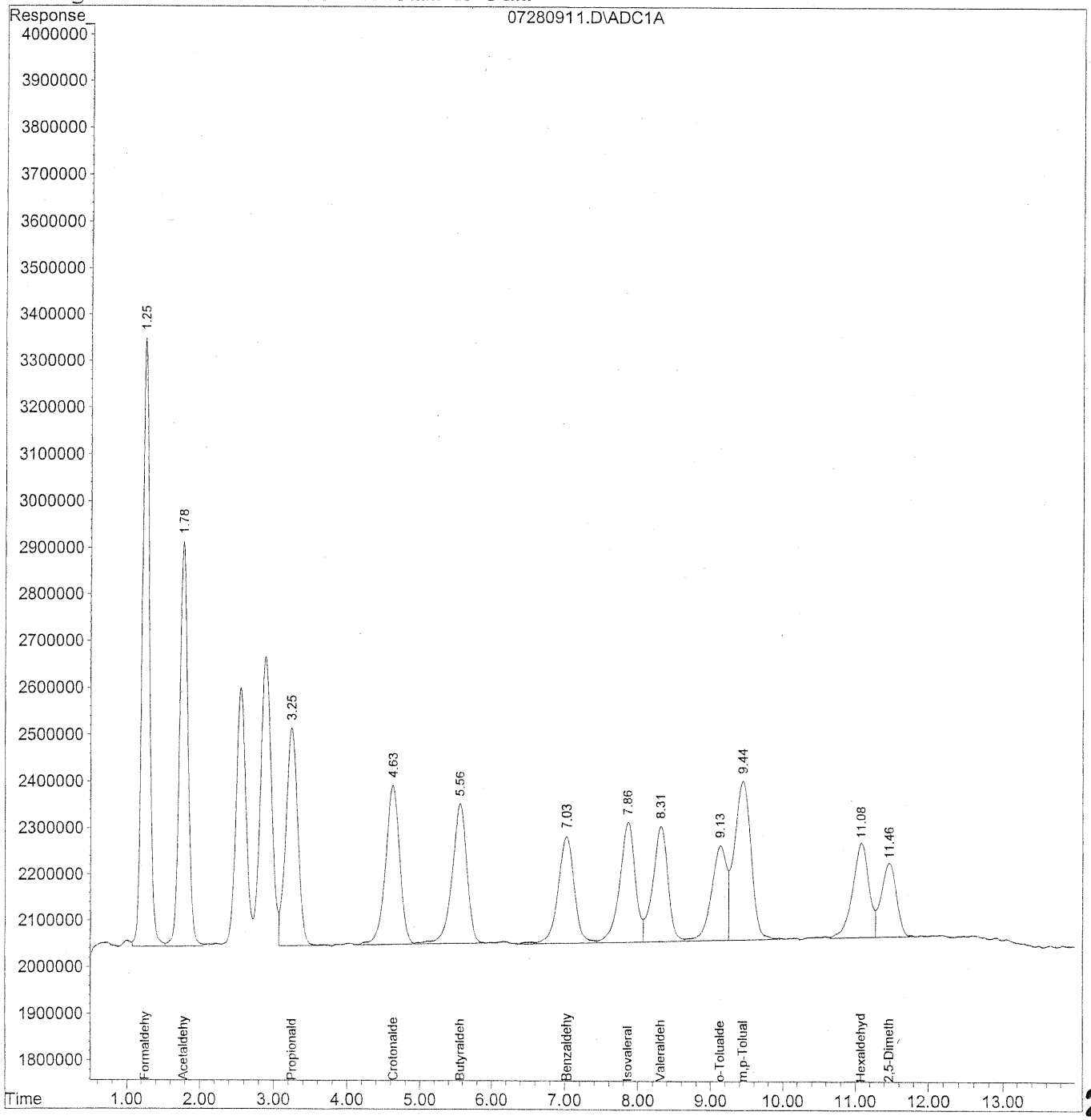
(f)=RT Delta > 1/2 Window (m)=manual int.

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



Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280911.D Vial: 11
Acq On : 28 Jul 2009 11:09 am Operator: HC
Sample : 500ng/ml TO11A Std S21-07270904 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:41 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

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

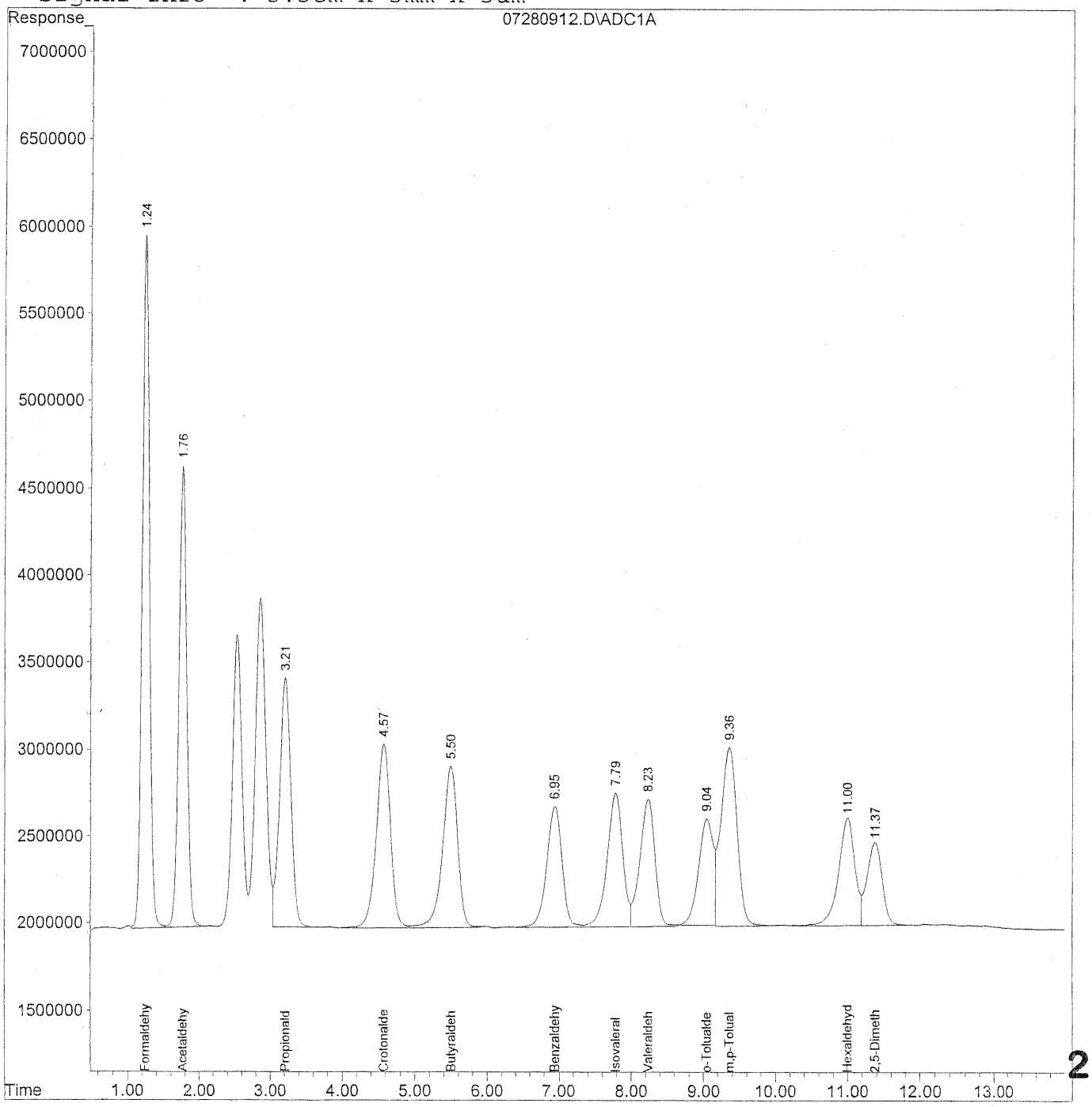
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.25	91399555	520.520	ng/ml
2)	Acetaldehyde	1.78	69908753	518.229	ng/ml
3)	Propionaldehyde	3.25	52190620	507.713	ng/ml
4)	Crotonaldehyde	4.63	46362546	419.361	ng/ml
5)	Butyraldehyde	5.56	43673214	542.691	ng/ml
6)	Benzaldehyde	7.03	34084716	540.400	ng/ml
7)	Isovaleraldehyde	7.87	39175205	441.901	ng/ml
8)	Valeraldehyde	8.31	36501988	439.281	ng/ml
9)	o-Tolualdehyde	9.13	30169058	560.032	ng/ml
10)	m,p-Tolualdehyde	9.44	54668231	1014.919	ng/ml
11)	Hexaldehyde	11.08	32179520	479.300	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.46	23309464	448.831	ng/ml

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

Data File : J:\LC01\DATA\TO11\2009_07\28\07280912.D Vial: 12
Acq On : 28 Jul 2009 11:24 am Operator:
Sample : 1500ng/ml TO11A Std S21-07270903 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:45 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

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



213

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

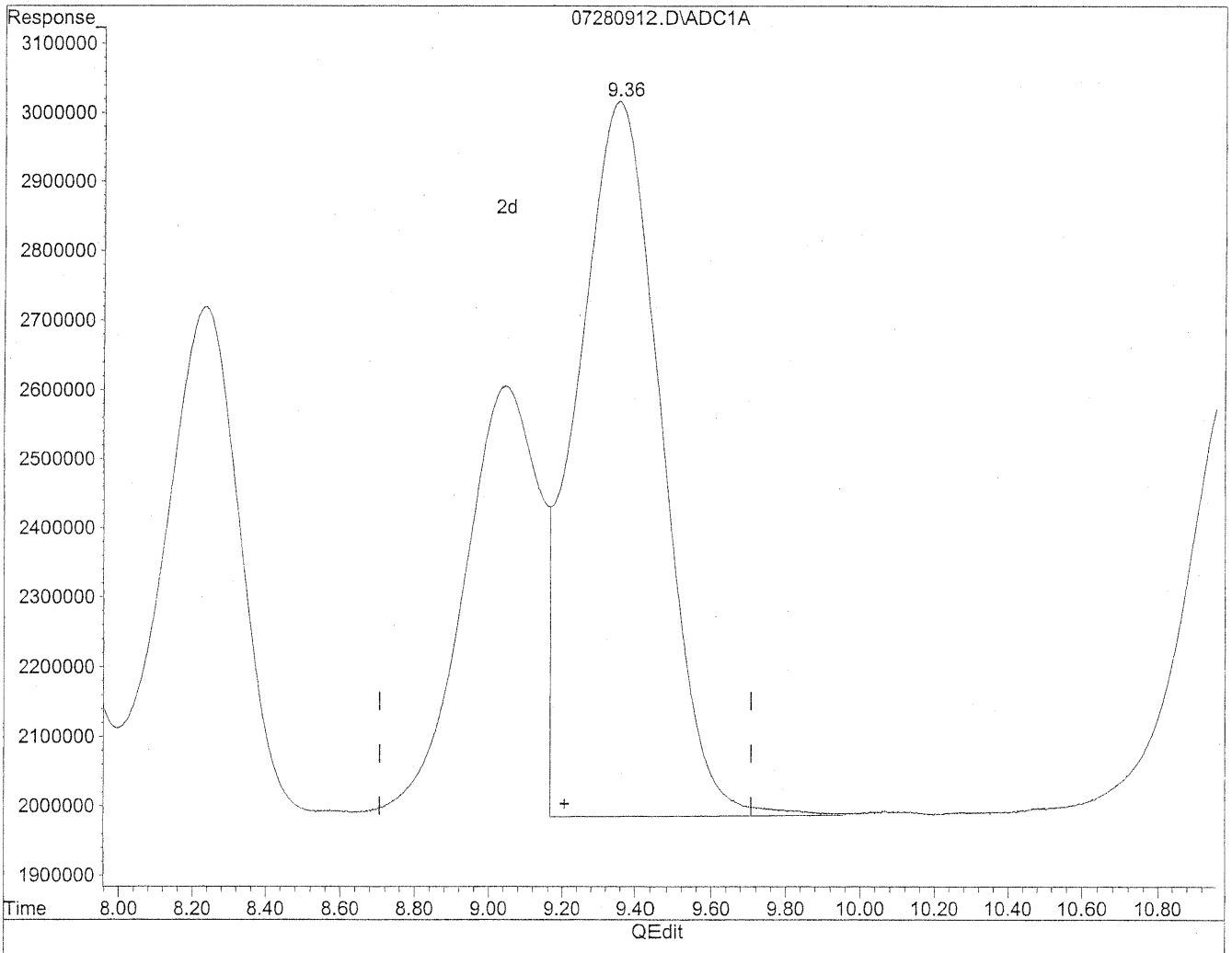
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.24	275380897	1568.292	ng/ml
2)	Acetaldehyde	1.76	209374751	1552.082	ng/ml
3)	Propionaldehyde	3.21	159030091	1547.054	ng/ml
4)	Crotonaldehyde	4.57	143227783	1295.530	ng/ml
5)	Butyraldehyde	5.50	134132687	1666.757	ng/ml
6)	Benzaldehyde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

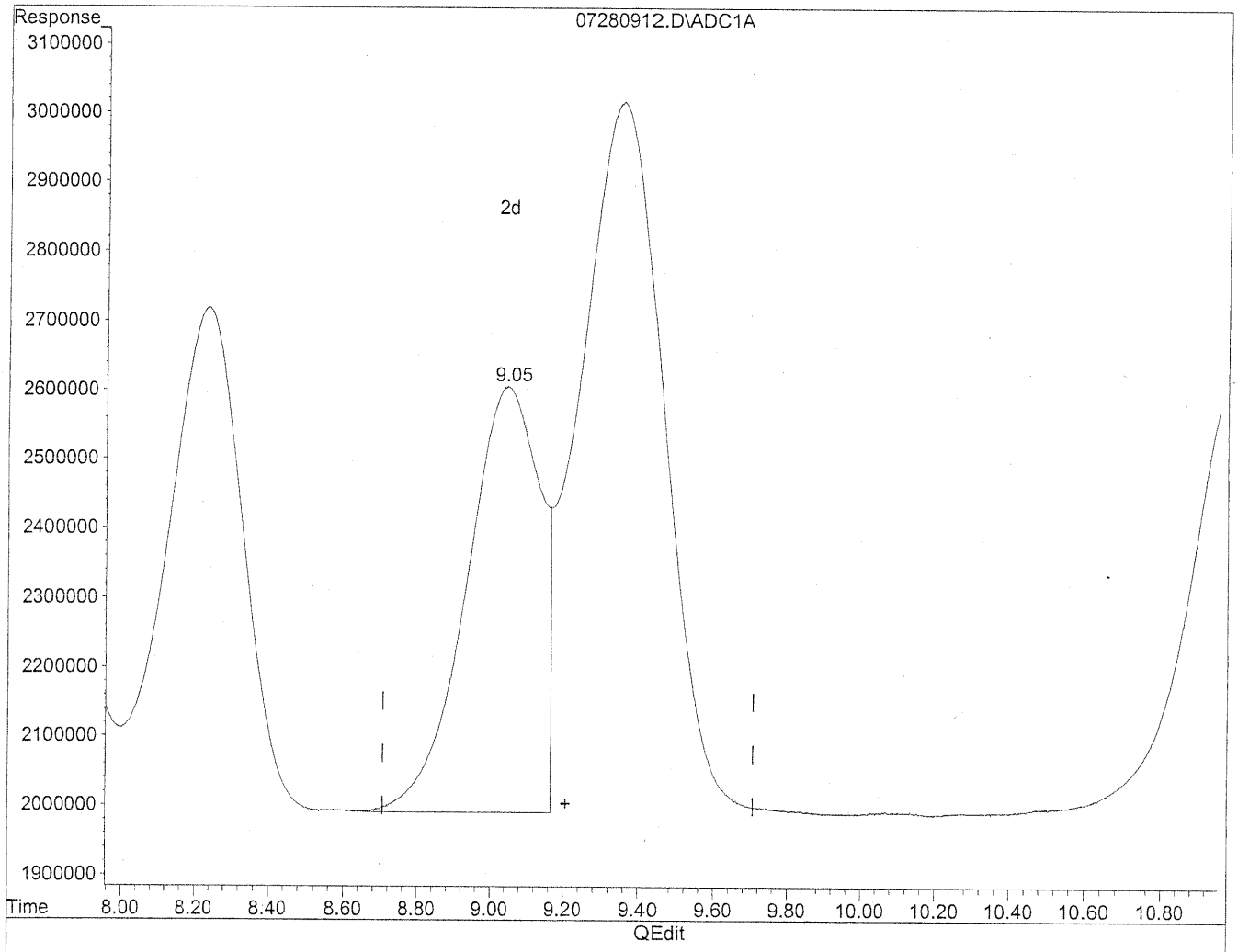
215

(+) = 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

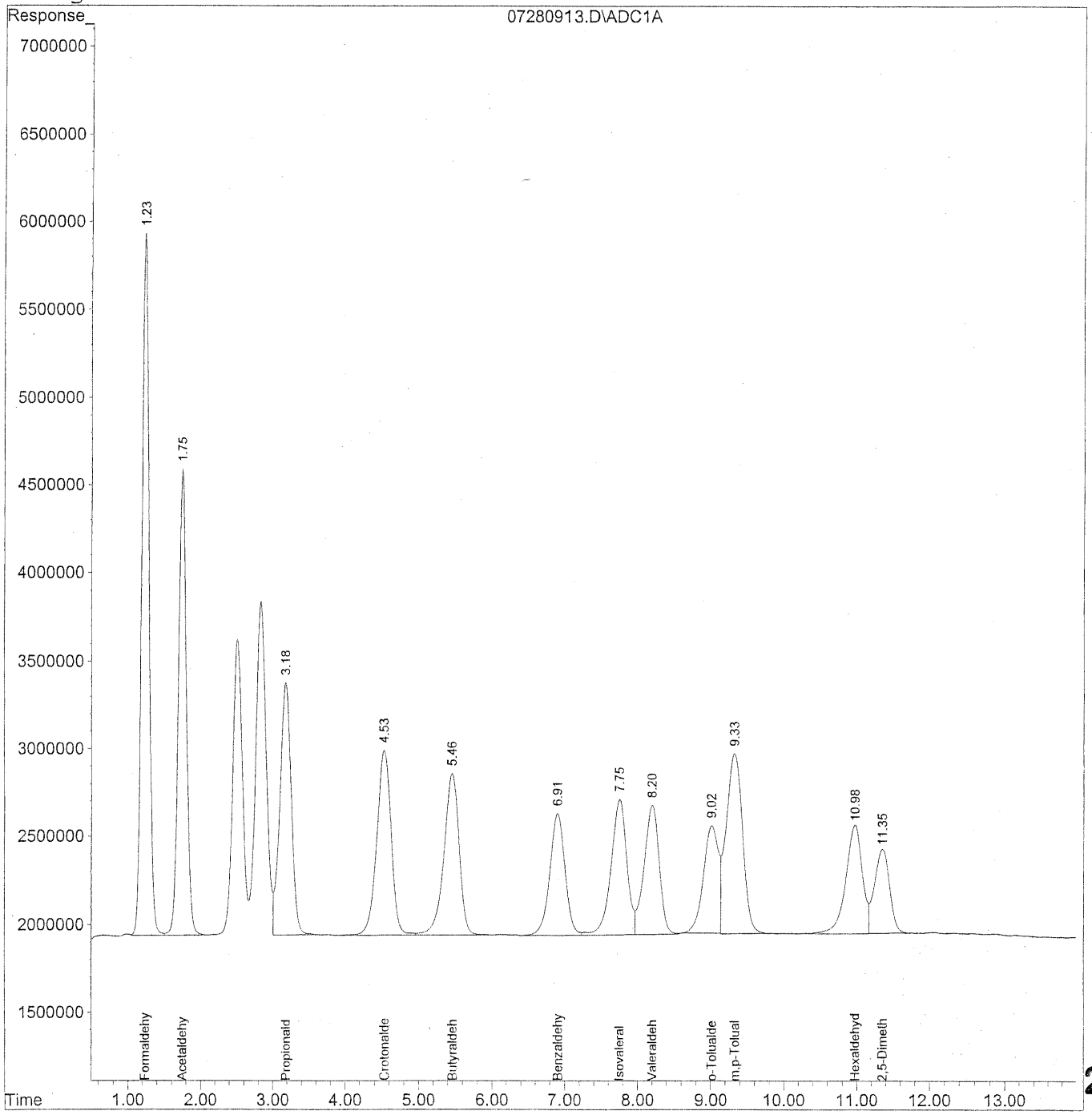
(+) = Expected Retention Time

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 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



217

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

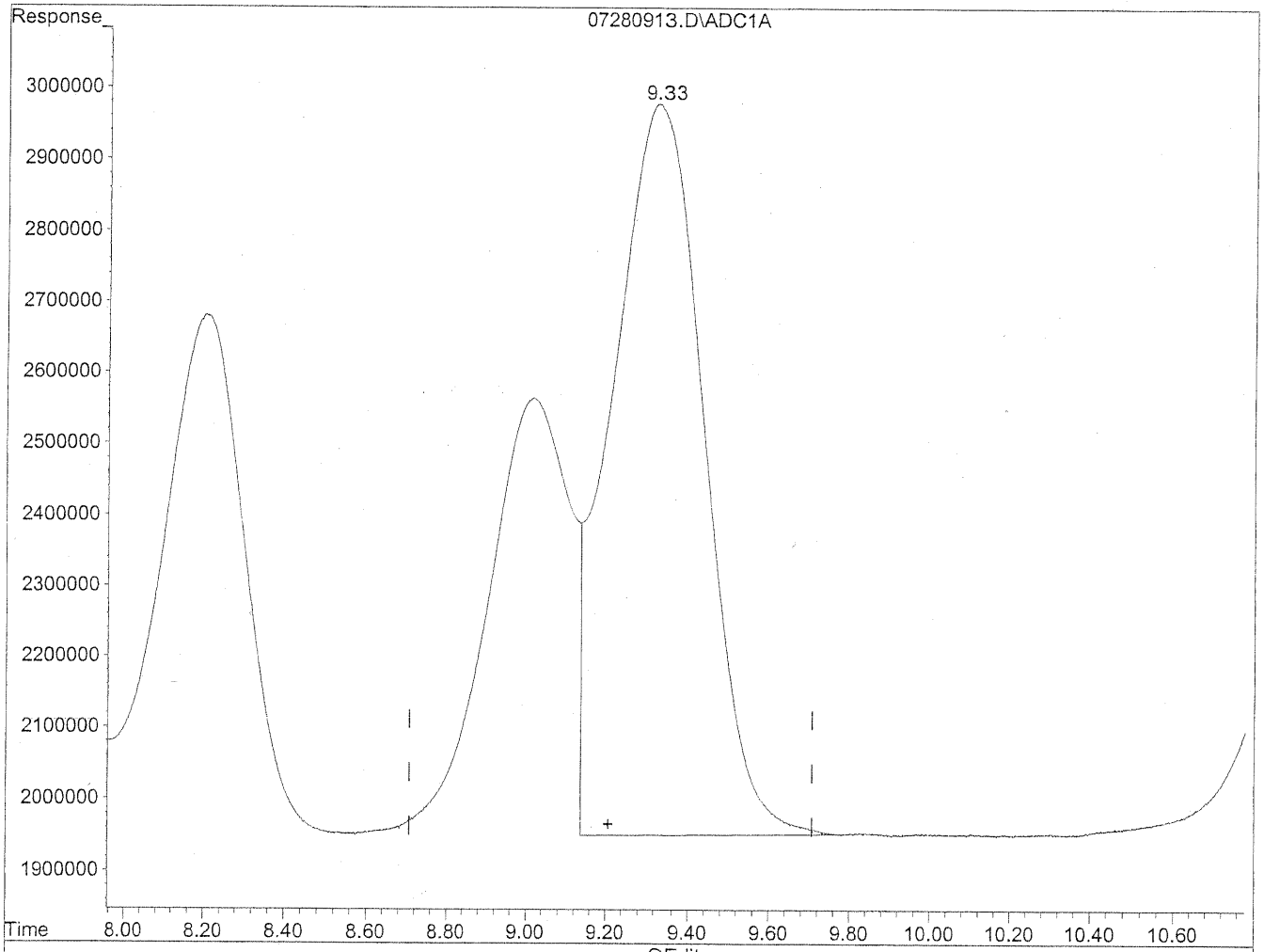
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.23	274724982	1564.557	ng/ml
2)	Acetaldehyde	1.75	209301649	1551.540	ng/ml
3)	Propionaldehyde	3.18	158919579	1545.979	ng/ml
4)	Crotonaldehyde	4.53	142112419	1285.442	ng/ml
5)	Butyraldehyde	5.46	132549734	1647.087	ng/ml
6)	Benzaldehyde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

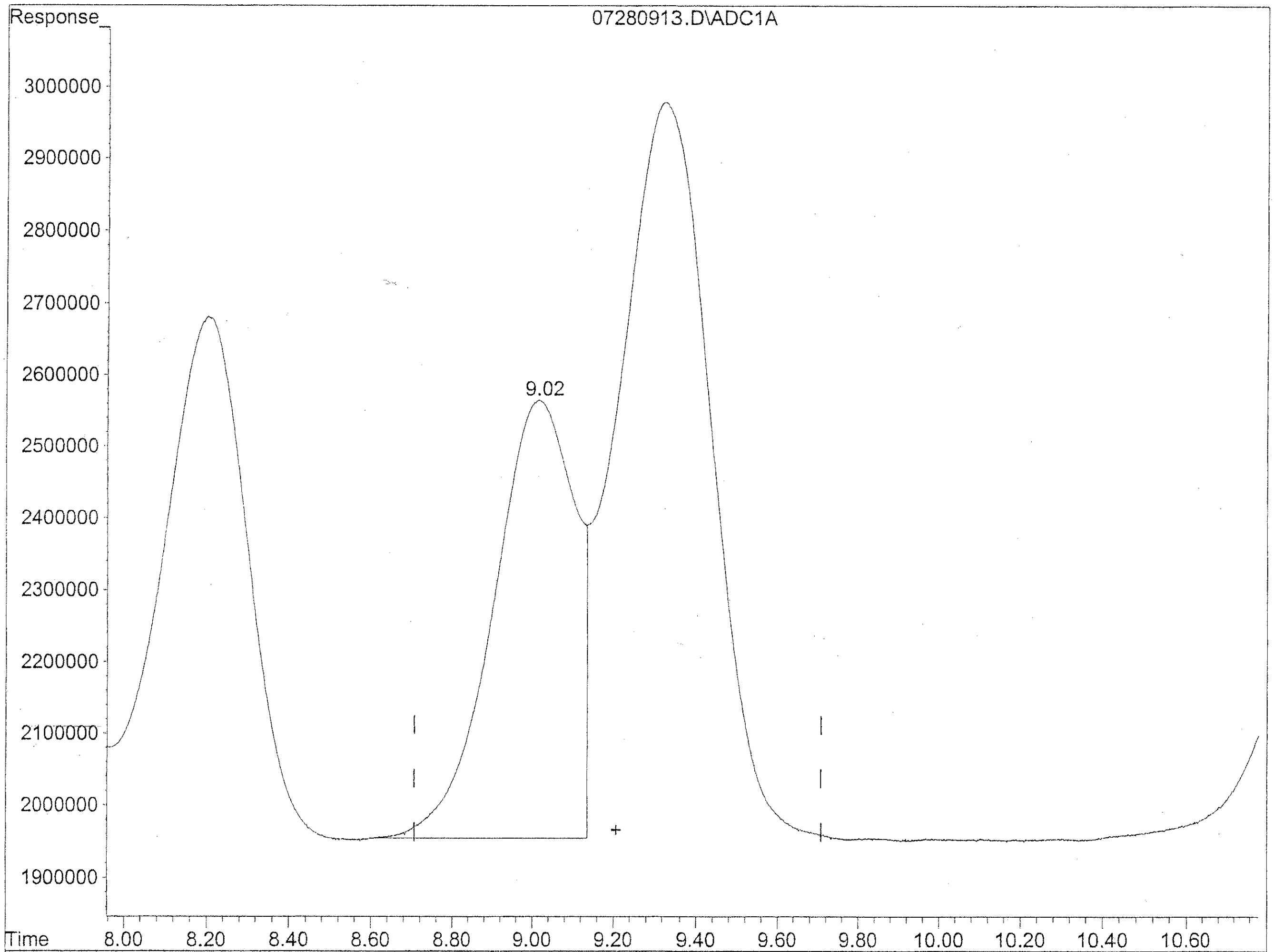
response 161094009

(+) = Expected Retention Time
07280913.D TO110709.M Tue Jul 28 14:47:05 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280913.D   Vial: 13
Acq On : 28 Jul 2009 11:39 am   Operator: HC
Sample : 1500ng/ml TO11A Std   Inst : LC 01
Misc :   Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:46 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



QEdit

(9) o-Tolualdehyde

9.02min 1595.318ng/ml m

response 85940120

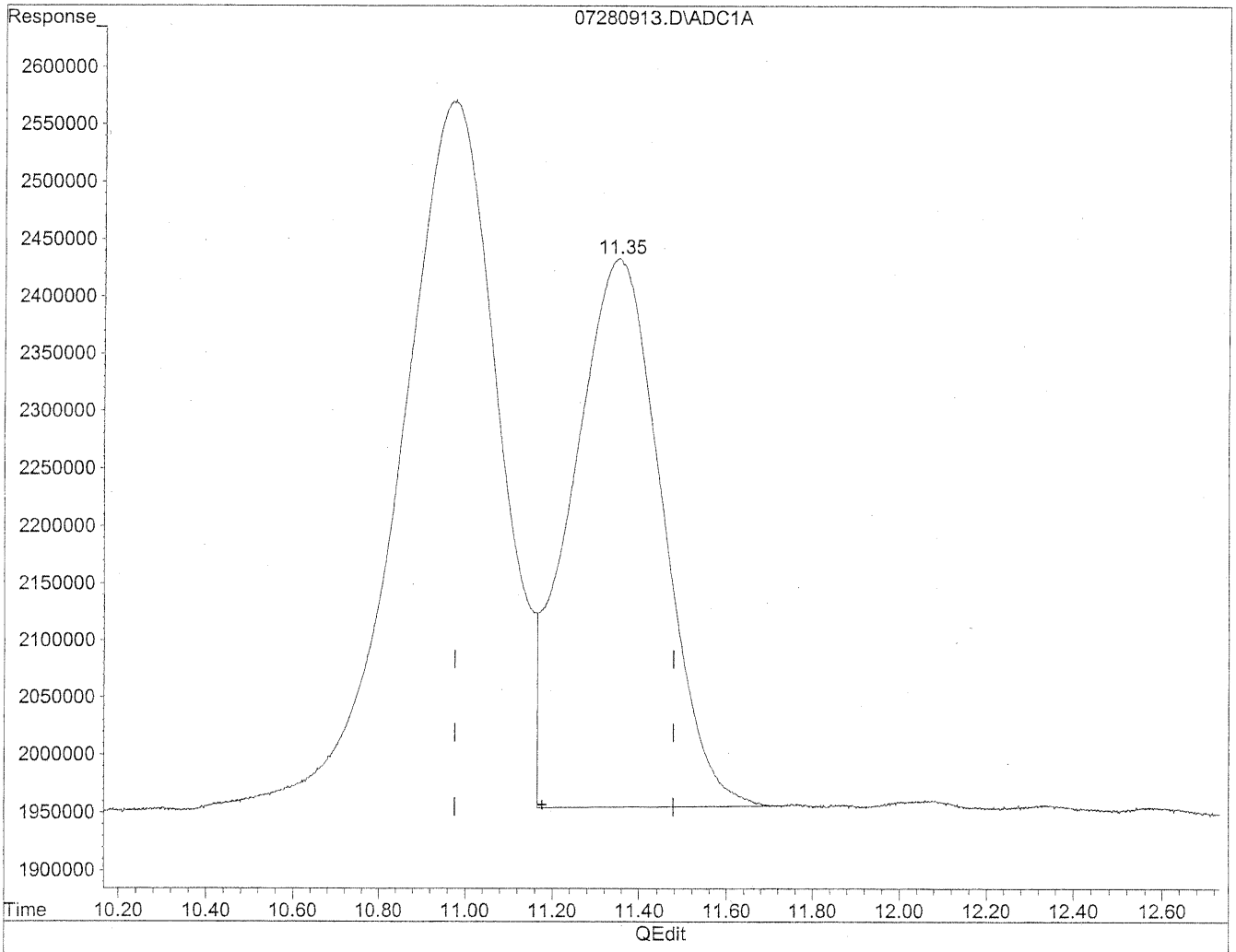
(+) = Expected Retention Time

07280913.D TO110709.M Tue Jul 28 14:47:19 2009

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280913.D Vial: 13
Acq On : 28 Jul 2009 11:39 am Operator: HC
Sample : 1500ng/ml TO11A Std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:46 19109 Quant Results File: TO110709.RES

Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration



(11) Hexaldehyde

11.35min 1025.842ng/ml

response 68873541

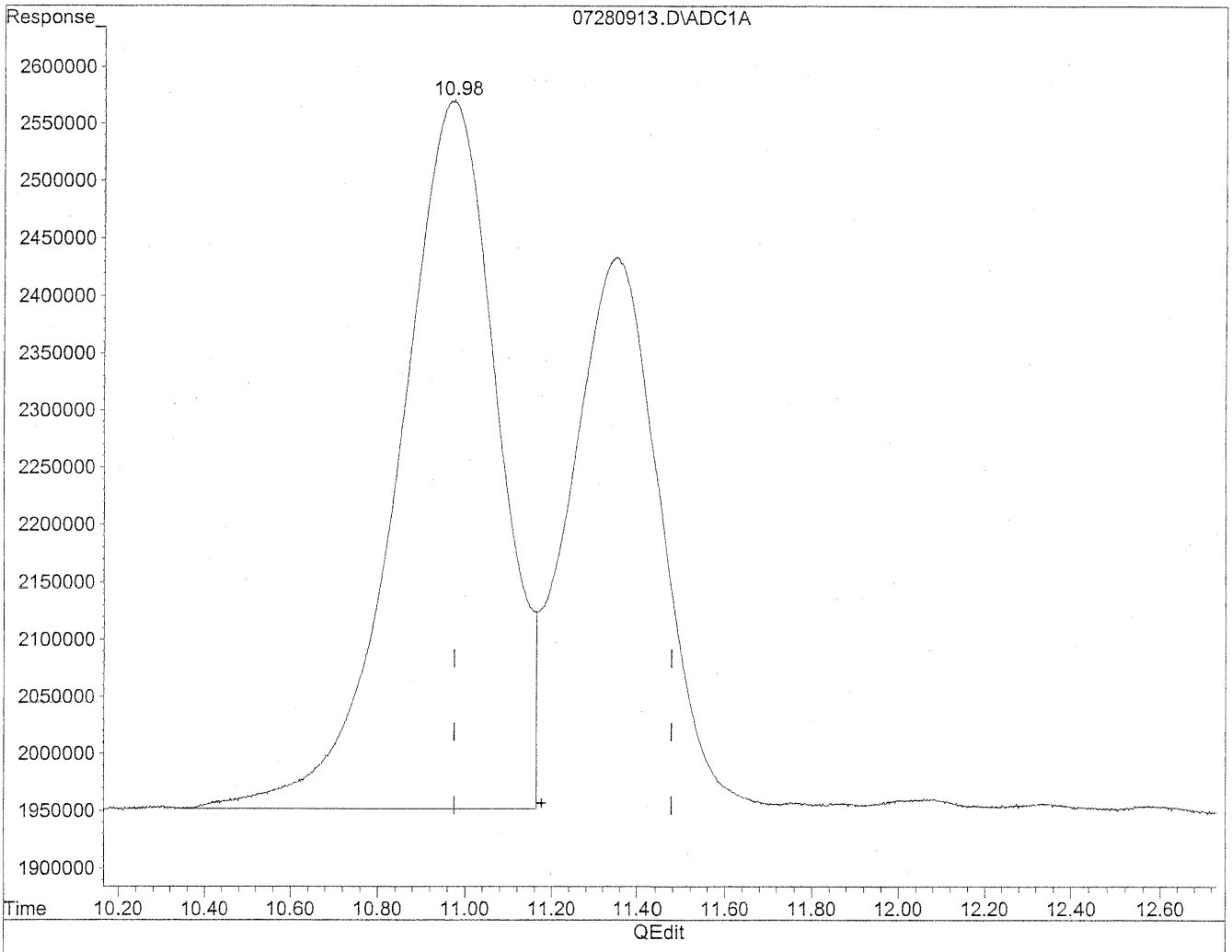
221

(+) = 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

10.98min 1461.011ng/ml m

response 98090122

KL 7/29/09

222

(+) = Expected Retention Time

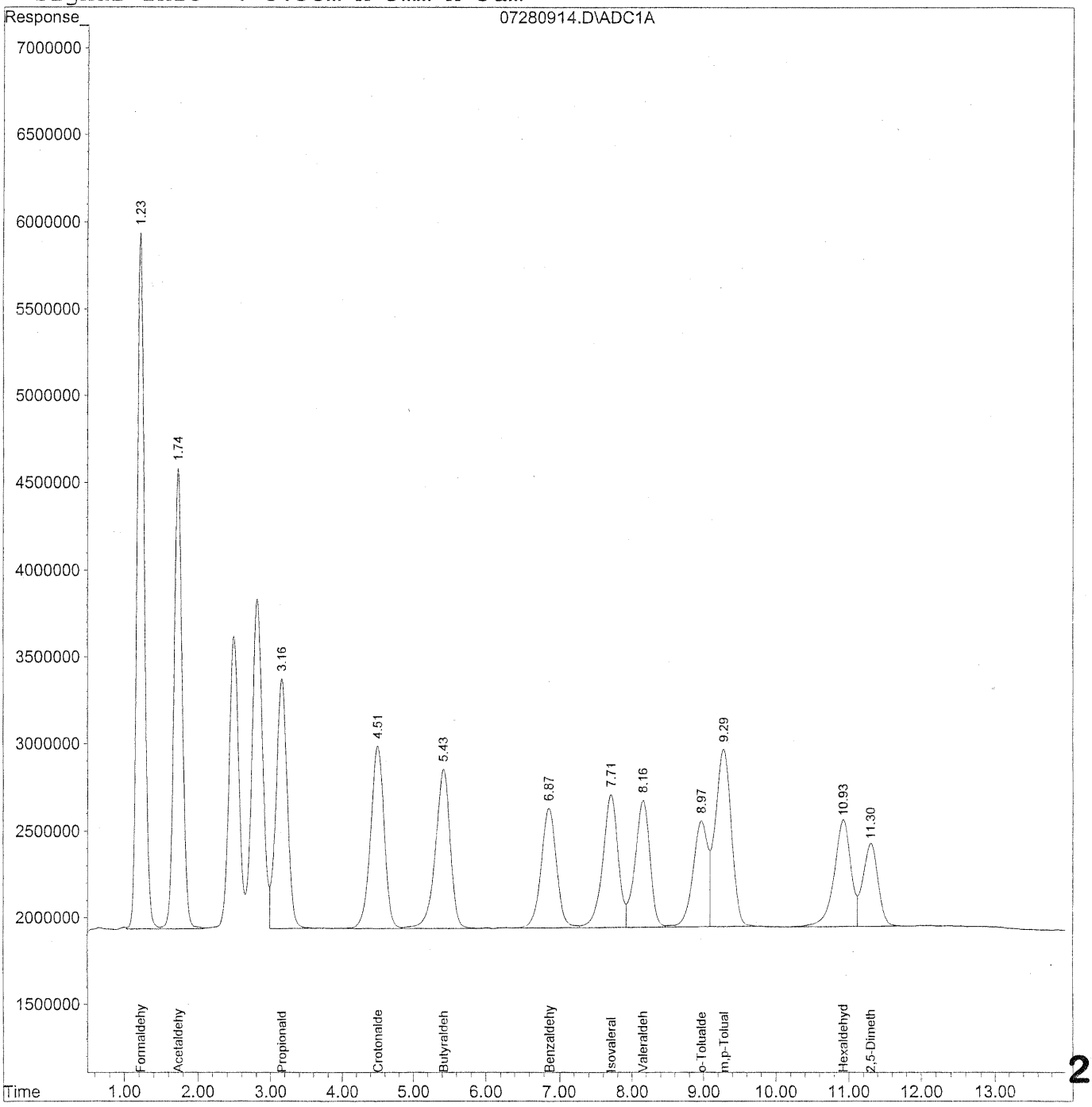
07280913.D TO110709.M Tue Jul 28 14:47:35 2009

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



223

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

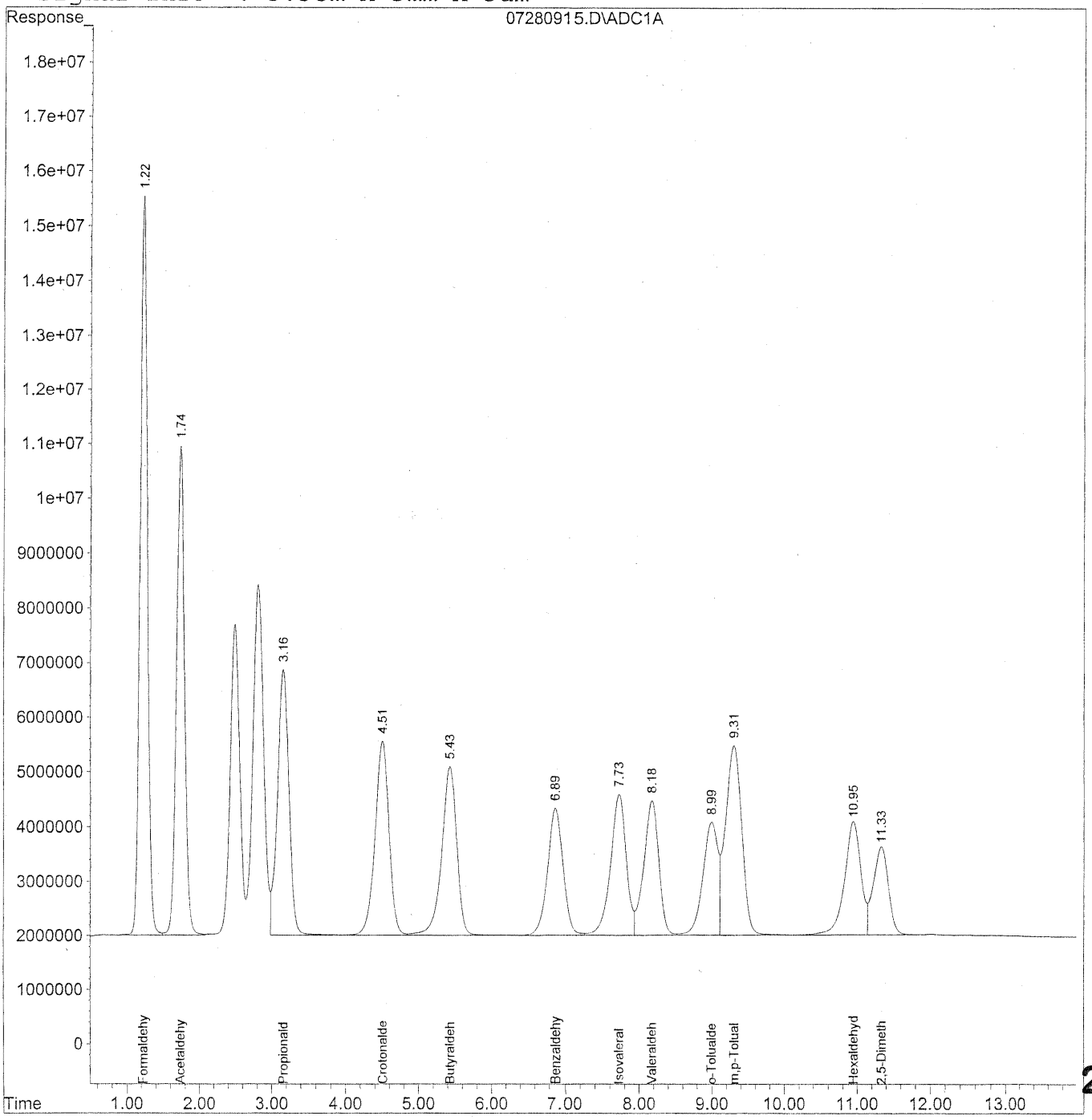
(f)=RT Delta > 1/2 Window (m)=manual int.

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



225

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280915.D Vial: 15
Acq On : 28 Jul 2009 12:09 pm Operator: HC
Sample : 5000ng/ml TO11A Std S21-07270902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:48 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

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

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	928364658	5175.655	ng/ml
2)	Acetaldehyde	1.74	706170560	5146.384	ng/ml
3)	Propionaldehyde	3.16	539067854	5164.924	ng/ml
4)	Crotonaldehyde	4.51	476268543	4639.802	ng/ml
5)	Butyraldehyde	5.43	446392739	5287.707	ng/ml
6)	Benzaldehyde	6.89	328286106	5108.361	ng/ml
7)	Isovaleraldehyde	7.73	388247386	4664.549	ng/ml
8)	Valeraldehyde	8.18	357832844	4632.141	ng/ml
9)	o-Tolualdehyde	8.99	298513860	5370.710	ng/ml
10)	m,p-Tolualdehyde	9.31	545640330	10145.539	ng/ml
11)	Hexaldehyde	10.95	332315493	5014.874	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.33	235692401	4817.453	ng/ml

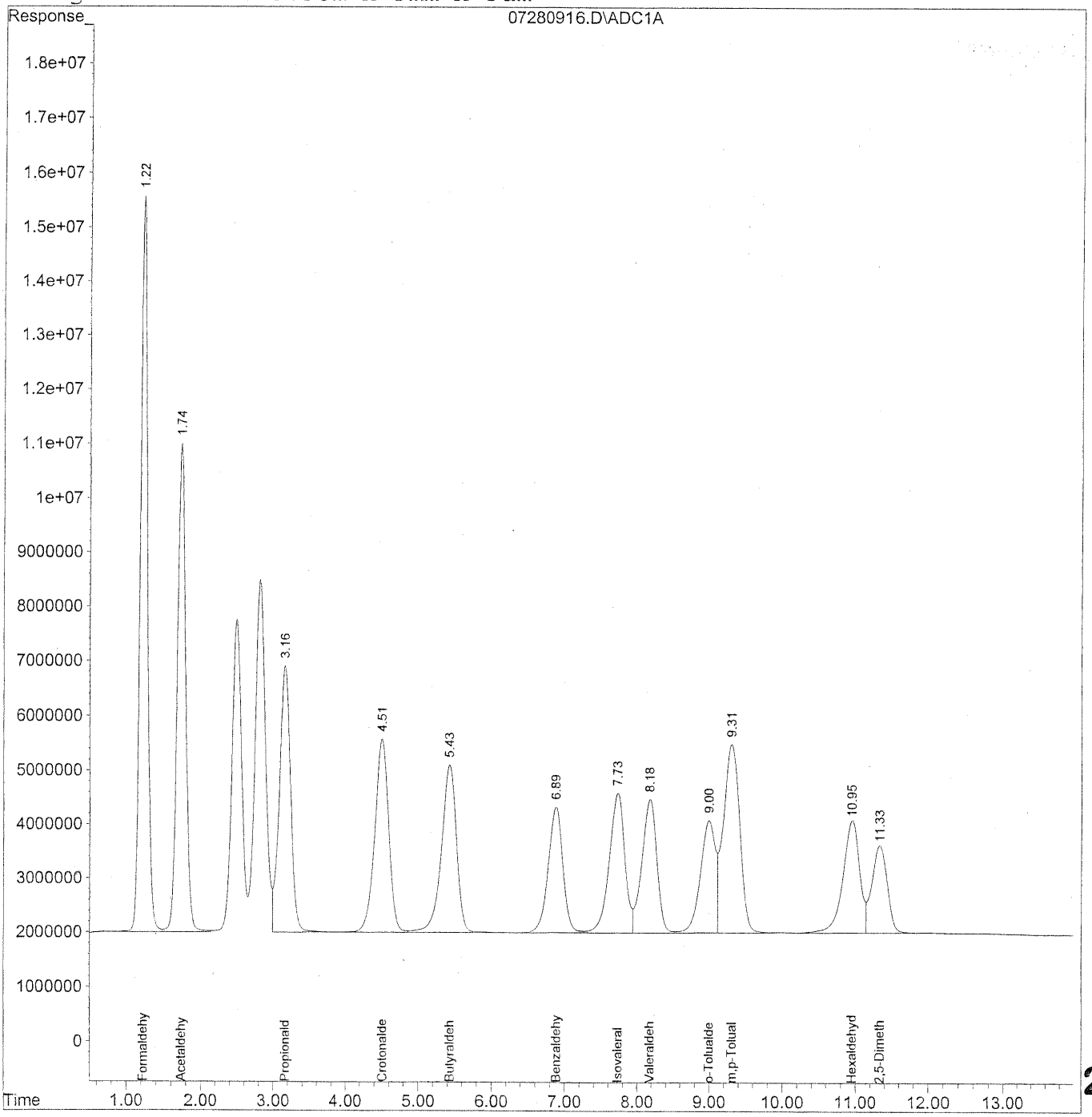
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280916.D Vial: 16
Acq On : 28 Jul 2009 12:24 pm Operator: HC
Sample : 5000ng/ml TO11A Std S21-07270902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:48 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

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



227

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

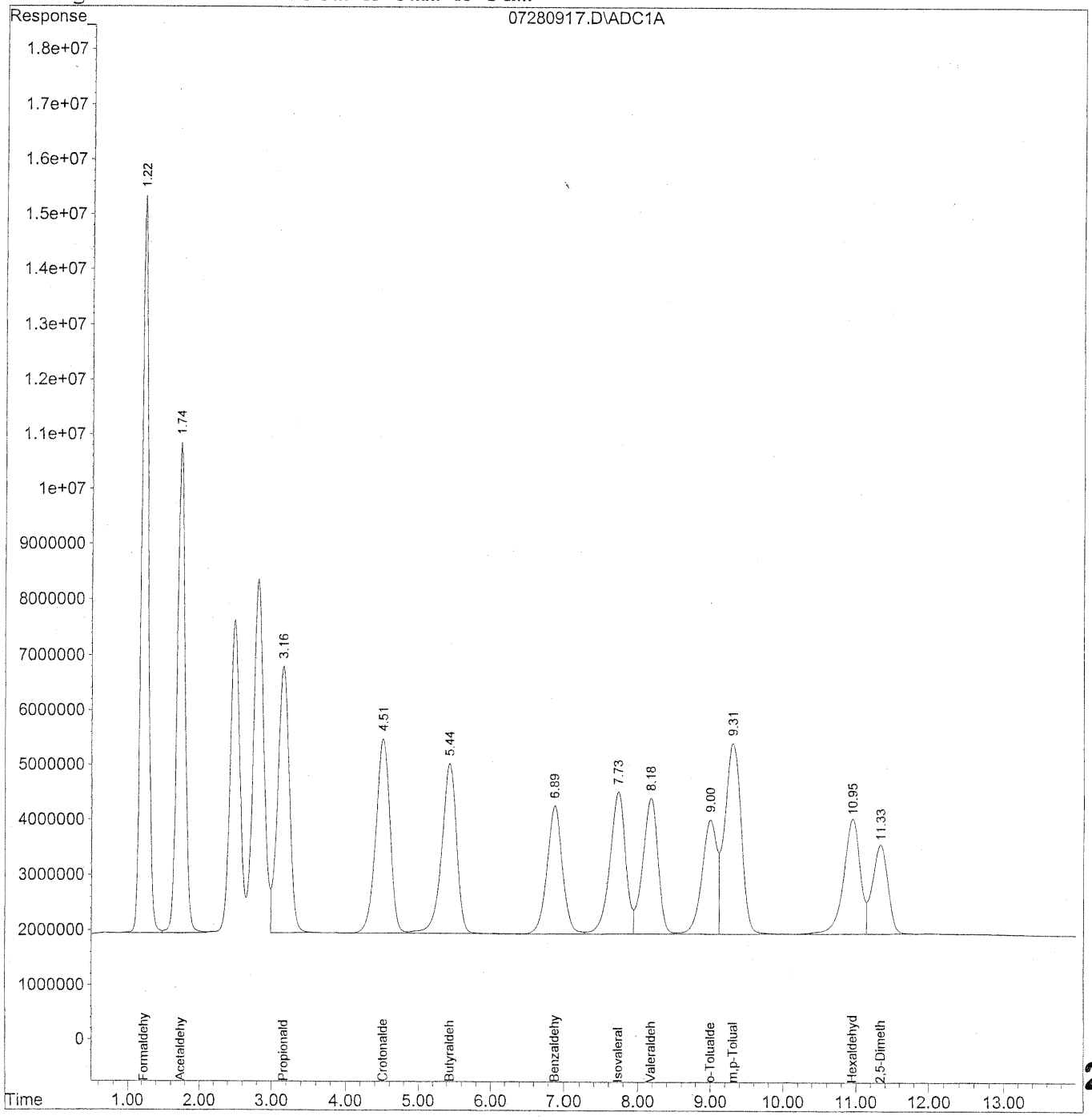
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280917.D Vial: 17
Acq On : 28 Jul 2009 12:39 pm Operator: HC
Sample : 5000ng/ml TO11A Std S21-07270902 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:49 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

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



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

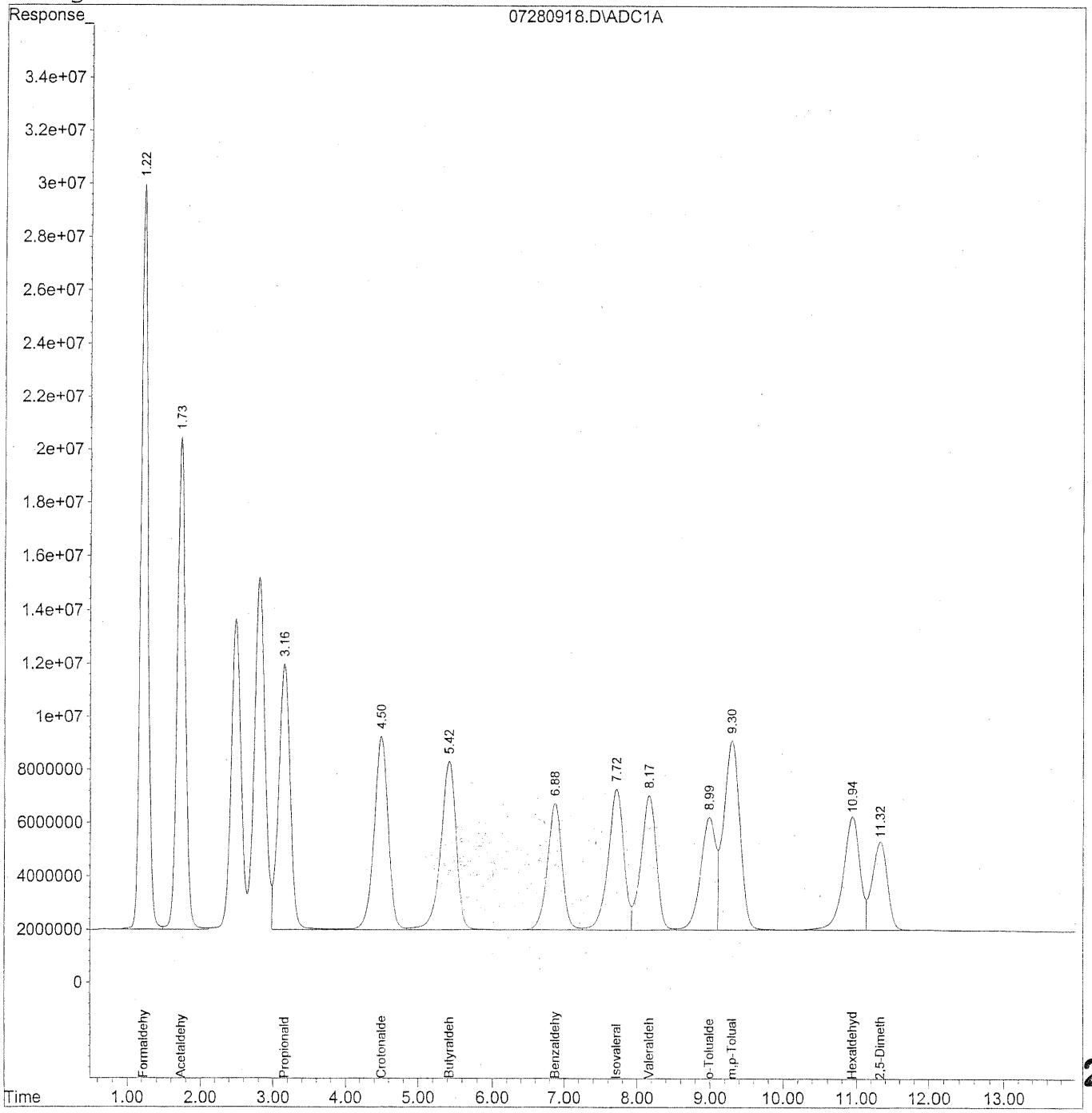
(f)=RT Delta > 1/2 Window (m)=manual int.

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



231

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

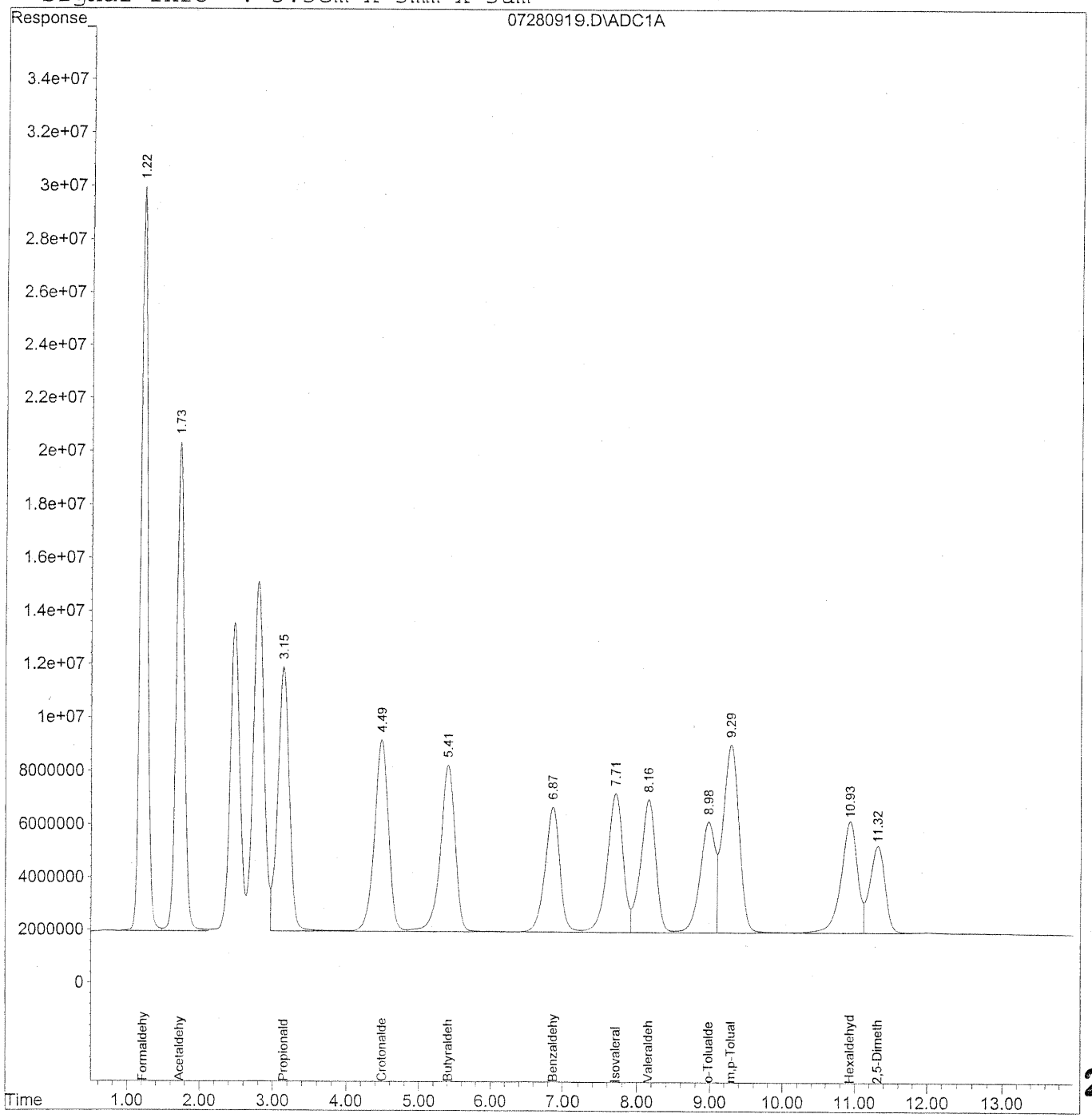
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

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



233

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280919.D Vial: 19
Acq On : 28 Jul 2009 1:09 pm Operator: HC
Sample : 10000ng/ml TO11A Std S21-07270901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:50 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

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

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	1905913073	10625.511	ng/ml
2)	Acetaldehyde	1.73	1446499891	10541.708	ng/ml
3)	Propionaldehyde	3.15	1098837646	10528.198	ng/ml
4)	Crotonaldehyde	4.49	971357788	9462.954	ng/ml
5)	Butyraldehyde	5.41	911328243	10795.060	ng/ml
6)	Benzaldehyde	6.87	669128969	10412.114	ng/ml
7)	Isovaleraldehyde	7.71	788026190	9467.640	ng/ml
8)	Valeraldehyde	8.16	729839210	9447.758	ng/ml
9)	o-Tolualdehyde	8.98	610326238	10980.681	ng/ml
10)	m,p-Tolualdehyde	9.29	1113209810	20698.824	ng/ml
11)	Hexaldehyde	10.93	681915785	10290.587	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.32	484763918	9908.370	ng/ml

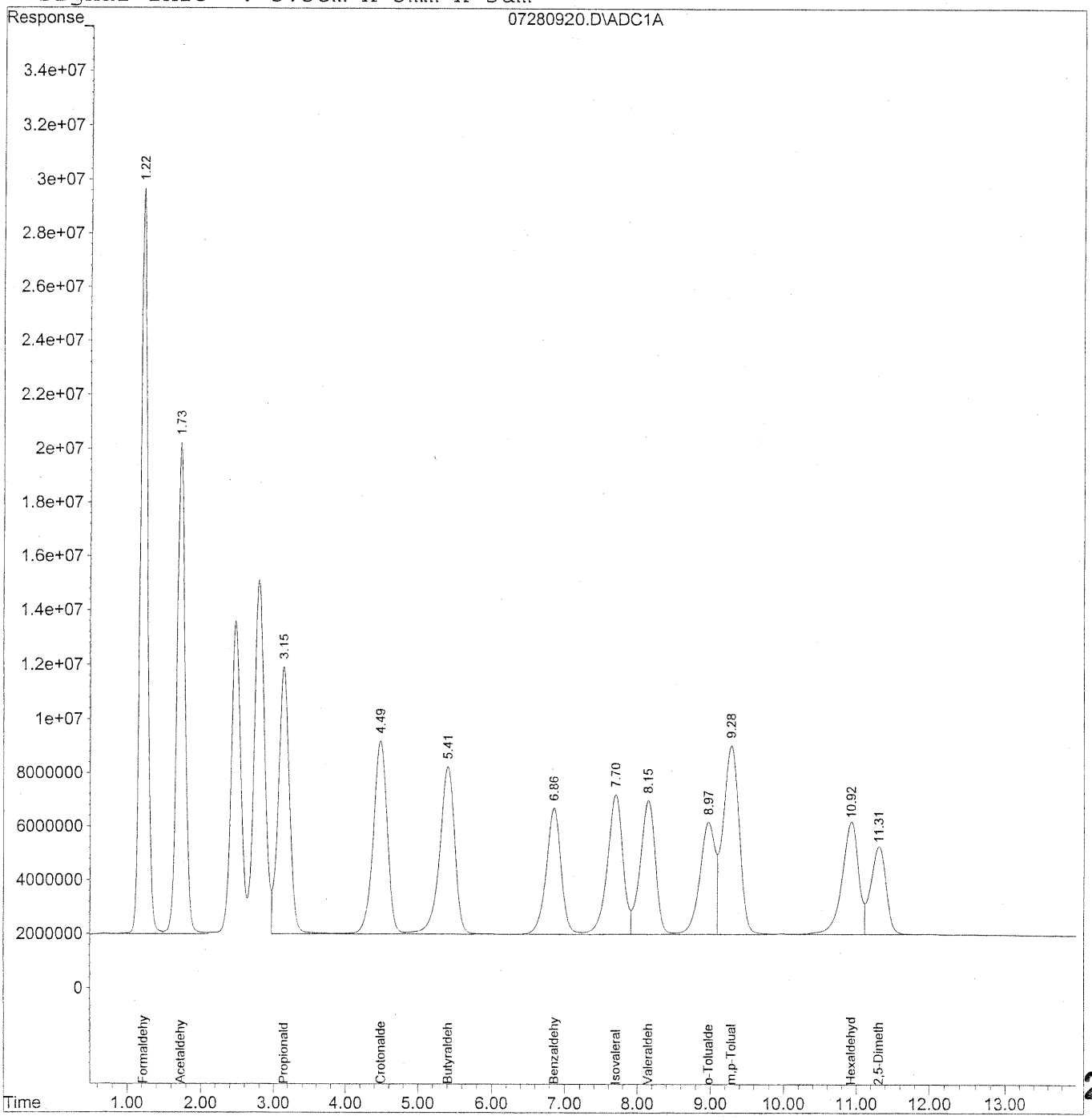
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280920.D Vial: 20
Acq On : 28 Jul 2009 1:25 pm Operator: HC
Sample : 10000ng/ml TO11A Std S21-07270901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:50 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

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



235

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280920.D Vial: 20
Acq On : 28 Jul 2009 1:25 pm Operator: HC
Sample : 10000ng/ml TO11A Std S21-07270901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 14:50 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 10:16:15 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

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

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	1875917434	10458.285	ng/ml
2)	Acetaldehyde	1.73	1425028469	10385.230	ng/ml
3)	Propionaldehyde	3.15	1089338811	10437.188	ng/ml
4)	Crotonaldehyde	4.48	963283335	9384.293	ng/ml
5)	Butyraldehyde	5.41	900561239	10667.520	ng/ml
6)	Benzaldehyde	6.86	662238443	10304.892	ng/ml
7)	Isovaleraldehyde	7.70	782256804	9398.325	ng/ml
8)	Valeraldehyde	8.15	722749626	9355.983	ng/ml
9)	o-Tolualdehyde	8.97	603256599	10853.487	ng/ml
10)	m,p-Tolualdehyde	9.29	1100384573	20460.354	ng/ml
11)	Hexaldehyde	10.92	670193360	10113.688	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.31	476113656	9731.563	ng/ml

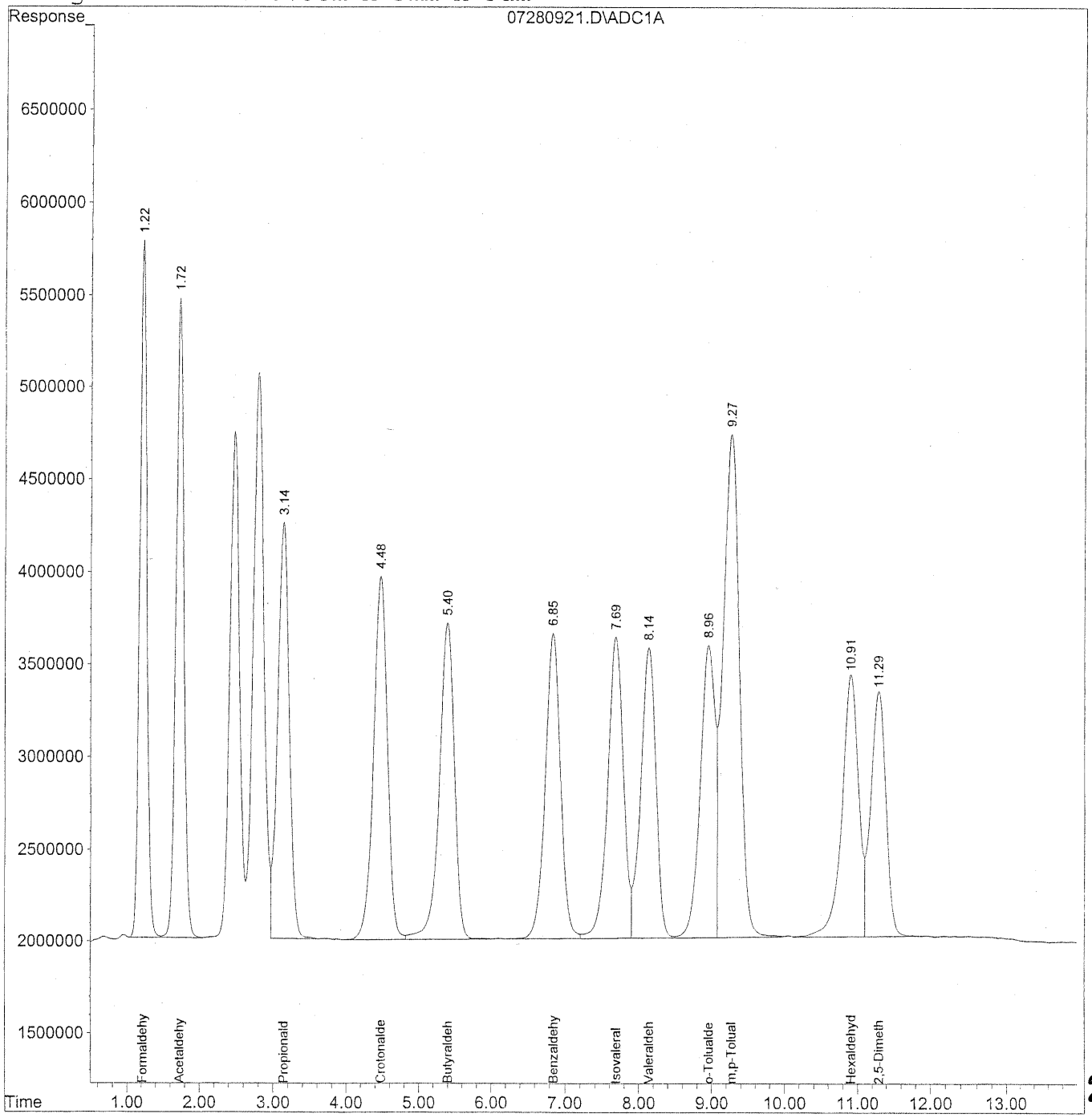
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_07\28\07280921.D Vial: 21
Acq On : 28 Jul 2009 1:40 pm Operator: HC
Sample : ~1500ng/ml TO11A Std ICV S21-07270907 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 17:22 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 15:29:52 2009
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

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



237

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_07\28\07280921.D Vial: 21
Acq On : 28 Jul 2009 1:40 pm Operator: HC
Sample : ~1500ng/ml TO11A Std ICV S21-07270907 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jul 28 17:22 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jul 28 15:29:52 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

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

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.22	257076667	1400.342	ng/ml
2)	Acetaldehyde	1.72	270257005	1927.330	ng/ml
3)	Propionaldehyde	3.14	246366252	2309.065	ng/ml
4)	Crotonaldehyde	4.48	262943470	2699.204	ng/ml
5)	Butyraldehyde	5.40	247400524	2800.672	ng/ml
6)	Benzaldehyde	6.85	233067402	3538.331	ng/ml
7)	Isovaleraldehyde	7.69	244473332	3002.720	ng/ml
8)	Valeraldehyde	8.14	226800810	3085.515	ng/ml
9)	o-Tolualdehyde	8.96	225349526	3863.990	ng/ml
10)	m,p-Tolualdehyde	9.27	428359795	7933.265	ng/ml
11)	Hexaldehyde	10.91	226495334	3363.271	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.29	193343187	3944.701	ng/ml

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

TO-11A Aldehyde-DNPH Stock Solution Standard S21-06300801

Source: AccuStandard Inc.
Catalog No: M-8315-R2-DNPH
Lot : B8060121
Solvent: ACN
Expiration Date: 6/12/11

HC 7/29/09

	MW	Aldehyde-DNPH MW*	Manufacturer Prepared Concentration as Aldehyde-DNPH (ug/mL)	Calculated Concentration as Aldehyde (ug/mL)	ICV S21-07270907 (nominal ng/mL)	ICV S21-07270907 (Actual, ng/mL)	% Diff
Formaldehyde	30.03	210.03	100	14.30	1430	1400.34	2.07%
Acetaldehyde	44.05	224.05	100.2	19.70	1970	1927.33	2.17%
Acetone	58.08	238.08	100.2	24.44	2444	not reported	
Acrolein	56.06	236.06	103.1	24.48	2448	not reported	
Propionaldehyde	58.08	238.08	100.2	24.44	2444	2309.07	5.52%
Crotonaldehyde	70.09	250.09	100.2	28.08	2808	2699.20	3.87%
Butyraldehyde	72.11	252.11	100	28.60	2860	2800.67	2.07%
Benzaldehyde	106.12	286.12	100	37.09	3709	3538.33	4.60%
Isovaleraldehyde	86.13	266.13	100.2	32.43	3243	3002.72	7.41%
Valeraldehyde	86.13	266.13	100.1	32.40	3240	3085.52	4.77%
o-Tolualdehyde	120.15	300.15	100.1	40.07	4007	3863.99	3.57%
m,p-Tolualdehyde	120.15	300.15	100.3	80.30	8030	7933.27	1.20%
Hexaldehyde	100.16	280.16	100.3	35.86	3586	3363.27	6.21%
2,5-Dimethylbenzaldehyde	134.18	314.18	100.3	42.84	4284	3944.70	7.92%

(* MW of DNPH is 198g/mol. The result of a nucleophilic reaction of aldehyde & DNPH is a hydrazone derivative with the loss of H20, 18g/mol)

CONTINUING CALIBRATION STANDARDS

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

HC 8/31/09

Printed : 8/31/09
Instrument : LC#1
Date Acquired : 8/24/09
Detector : UV-VIS 360
Sample Amount : 5ul
Analyst : HC
Client & PAI Job# : EH&E P0902912

SAMPLE RESULT SUMMARY

Sample Information	MDL	CCV 1500ng/ml S21-08240901	% Diff	ACN lot Blk CY023	MB front lot 5855/5994 1.0ml	MB back lot 5855/5994 1.0ml	P0902912-001 back 1.0ml	P0902912-002 back 1.0ml	P0902912-003 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	102.00	101.50	108.68
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	1374.8	8.3%	ND	ND	ND	148.529	ND	ND
Acetaldehyde	100.00	1358.8	9.4%	ND	ND	ND	556.821 BT	ND	567.946 BT
Propionaldehyde	100.00	1346.4	10.2%	ND	ND	ND	ND	ND	ND
Crotonaldehyde	100.00	1316.3	12.2%	ND	ND	ND	ND	ND	ND
Butyraldehyde	100.00	1363.3	9.1%	ND	ND	ND	ND	ND	ND
Benzaldehyde	100.00	1352.4	9.8%	ND	ND	ND	ND	ND	ND
Isovaleraldehyde	100.00	1335.8	10.9%	ND	ND	ND	ND	ND	ND
Valeraldehyde	100.00	1290.7	14.0%	ND	ND	ND	ND	ND	ND
o-Tolualdehyde	100.00	1370.0	8.7%	ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde	200.00	2685.2	10.5%	ND	ND	ND	ND	ND	ND
Hexaldehyde	100.00	1377.1	8.2%	ND	ND	ND	ND	ND	ND
2,5-Dimethylbenzaldehyde	100.00	1277.4	14.8%	ND	ND	ND	ND	ND	ND

ug/m3		ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3
Formaldehyde		NA	NA	NA	1.456	ND	ND
Acetaldehyde		NA	NA	NA	5.459	ND	5.226
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	1.186	ND	ND
Acetaldehyde		NA	NA	NA	3.031	ND	2.902
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 : 8/24/09
Detector : UV-VIS 360 Sample Amount : 5ul
Analyst : HC Client & PAI Job# : EH&E P0902912

SAMPLE RESULT SUMMARY

Sample Information	MDL	P0902912-004 back 1.0ml	P0902912-005 back 1.0ml	CCV 1500ng/ml S21-08240901	% Diff	CCV 1500ng/ml S21-08240901	% Diff	ACN blk lot CY023
Dilution	1.0	1.0	1.0	1.0		1.0		1.0
Sample Volume (L)	NA	102.96	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	1422.632	5.2%	1383.560	7.8%	ND
Acetaldehyde	100.00	390.720	ND	1406.872	6.2%	1383.308	7.8%	ND
Propionaldehyde	100.00	ND	ND	1390.934	7.3%	1368.198	8.8%	ND
Crotonaldehyde	100.00	ND	ND	1360.678	9.3%	1344.253	10.4%	ND
Butyraldehyde	100.00	ND	ND	1406.095	6.3%	1396.275	6.9%	ND
Benzaldehyde	100.00	ND	ND	1401.372	6.6%	1370.317	8.6%	ND
Isovaleraldehyde	100.00	ND	ND	1405.973	6.3%	1374.858	8.3%	ND
Valeraldehyde	100.00	ND	ND	1365.035	9.0%	1340.975	10.6%	ND
o-Tolualdehyde	100.00	ND	ND	1433.568	4.4%	1412.560	5.8%	ND
m,p-Tolualdehyde	200.00	ND	ND	2804.480	6.5%	2769.042	7.7%	ND
Hexaldehyde	100.00	ND	ND	1424.591	5.0%	1410.338	6.0%	ND
2,5-Dimethylbenzaldehyde	100.00	ND	ND	1275.502	15.0%	1297.339	13.5%	ND

	ug/m3	ug/m3		ug/m3
Formaldehyde	ND	ND		ND
Acetaldehyde	3.795	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
Formaldehyde	ND	ND		ND
Acetaldehyde	2.107	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

Printed : 8/31/09
Instrument : LC#1 Date Acquired : 8/24/09
Detector : UV-VIS 360 Sample Amount : 5ul
Analyst : HC Client & PAI Job# : EH&E P0902912

Sample Information	MDL	MB front lot 5855/5994 1.0ml	MB back lot 5855/5994 1.0ml	P0902912-001 front 1.0ml	P0902912-002 front 1.0ml	P0902912-003 front 1.0ml	P0902912-004 front 1.0ml	P0902912-005 front 1.0ml
Dilution	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
Sample Volume (L)	NA			102.00	101.50	108.68	102.96	0.00
Final Vol.(ml)	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0

	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample	ng/sample
Formaldehyde	100.00	ND	ND	5429.098	328.646	5801.044	5040.015	ND
Acetaldehyde	100.00	ND	ND	5466.173	ND	4941.131	4525.216	ND
Propionaldehyde	100.00	ND	ND	570.110	ND	529.272	473.759	ND
Crotonaldehyde	100.00	ND	ND	ND	ND	ND	ND	ND
Butyraldehyde	100.00	ND	ND	587.491	ND	489.857	481.216	ND
Benzaldehyde	100.00	ND	ND	396.650	ND	675.535	587.236	ND
Isovaleraldehyde	100.00	ND	ND	228.362	ND	208.097	190.687	ND
Valeraldehyde	100.00	ND	ND	1468.104	ND	1427.381	1312.042	ND
o-Tolualdehyde	100.00	ND	ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde	200.00	ND	ND	ND	ND	ND	ND	ND
Hexaldehyde	100.00	ND	ND	6451.295	ND	6264.929	5753.058	ND
2,5-Dimethylbenzaldehyde	100.00	ND	ND	ND	ND	ND	ND	ND

	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3
Formaldehyde		ND	ND	53.226	3.238	53.377	48.951	ND
Acetaldehyde		ND	ND	53.590	ND	45.465	43.951	ND
Propionaldehyde		ND	ND	5.589	ND	4.870	4.601	ND
Crotonaldehyde		ND	ND	ND	ND	ND	ND	ND
Butyraldehyde		ND	ND	5.760	ND	4.507	4.674	ND
Benzaldehyde		ND	ND	3.889	ND	6.216	5.704	ND
Isovaleraldehyde		ND	ND	2.239	ND	1.915	1.852	ND
Valeraldehyde		ND	ND	14.393	ND	13.134	12.743	ND
o-Tolualdehyde		ND	ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde		ND	ND	ND	ND	ND	ND	ND
Hexaldehyde		ND	ND	63.248	ND	57.646	55.877	ND
2,5-Dimethylbenzaldehyde		ND	ND	ND	ND	ND	ND	ND

	ppb	ppb	ppb	ppb	ppb	ppb	ppb	ppb
Formaldehyde		ND	ND	43.354	2.637	43.477	39.872	ND
Acetaldehyde		ND	ND	29.757	ND	25.246	24.405	ND
Propionaldehyde		ND	ND	2.354	ND	2.051	1.938	ND
Crotonaldehyde		ND	ND	ND	ND	ND	ND	ND
Butyraldehyde		ND	ND	1.954	ND	1.529	1.585	ND
Benzaldehyde		ND	ND	0.896	ND	1.433	1.315	ND
Isovaleraldehyde		ND	ND	0.636	ND	0.544	0.526	ND
Valeraldehyde		ND	ND	4.088	ND	3.730	3.619	ND
o-Tolualdehyde		ND	ND	ND	ND	ND	ND	ND
m,p-Tolualdehyde		ND	ND	ND	ND	ND	ND	ND
Hexaldehyde		ND	ND	15.446	ND	14.078	13.646	ND
2,5-Dimethylbenzaldehyde		ND	ND	ND	ND	ND	ND	ND

COLUMBIA ANALYTICAL SERVICES

TO11A Aldehyde & Ketone DNPH Analysis by HPLC

		Printed :	8/31/09
Instrument :	LC#1	Date Acquired	8/24/09
Detector :	UV-VIS 360	Sample Amoun	5ul
Analyst :	HC	Client & PAI Jo	EH&E P0902912

Sample Information	MDL	CCV 1500ng/ml S21-08240901	% Diff
Dilution	1.0	1.0	
Sample Volume (L)	NA		
Final Vol.(ml)	1.0	1.0	

	ng/sample	ng/sample	
Formaldehyde	100.00	1364.982	9.0%
Acetaldehyde	100.00	1360.629	9.3%
Propionaldehyde	100.00	1360.217	9.3%
Crotonaldehyde	100.00	1320.310	12.0%
Butyraldehyde	100.00	1363.354	9.1%
Benzaldehyde	100.00	1345.139	10.3%
Isovaleraldehyde	100.00	1327.393	11.5%
Valeraldehyde	100.00	1281.818	14.5%
o-Tolualdehyde	100.00	1341.338	10.6%
m,p-Tolualdehyde	200.00	2686.989	10.4%
Hexaldehyde	100.00	1333.370	11.1%
2,5-Dimethylbenzaldehyde	100.00	1278.602	14.8%

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

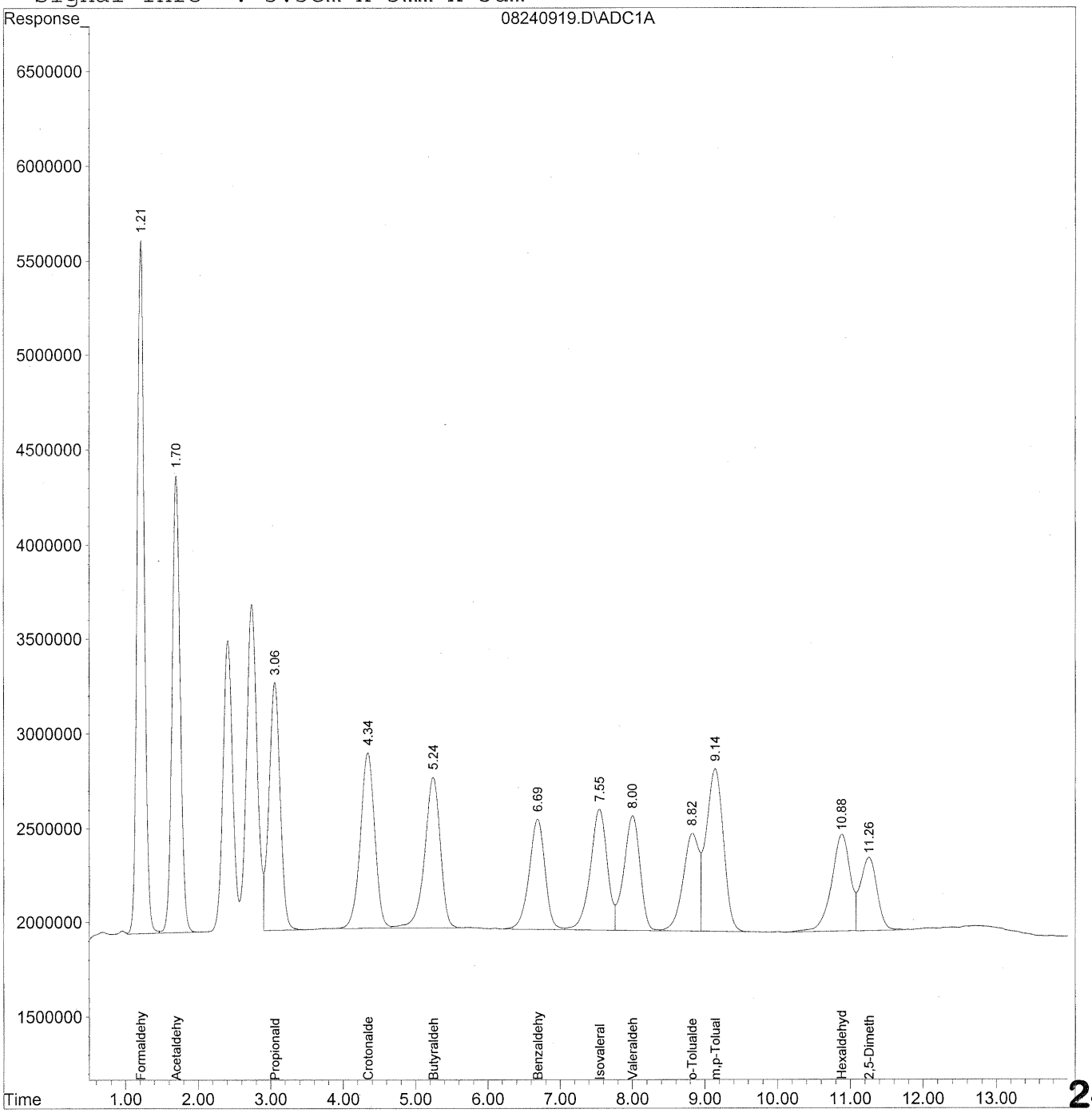
	ppb	
Formaldehyde		
Acetaldehyde		
Propionaldehyde		
Crotonaldehyde		
Butyraldehyde		
Benzaldehyde		
Isovaleraldehyde		
Valeraldehyde		
o-Tolualdehyde		
m,p-Tolualdehyde		
Hexaldehyde		
2,5-Dimethylbenzaldehyde		

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240919.D Vial: 17
Acq On : 24 Aug 2009 5:01 pm Operator: HC
Sample : CCV 1500ng/ml S21-08240901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 11:34 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 Aug 25 10:21:57 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\24\08240919.D Vial: 17
Acq On : 24 Aug 2009 5:01 pm Operator: HC
Sample : CCV 1500ng/ml S21-08240901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 11:34 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 Aug 25 10:21:57 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

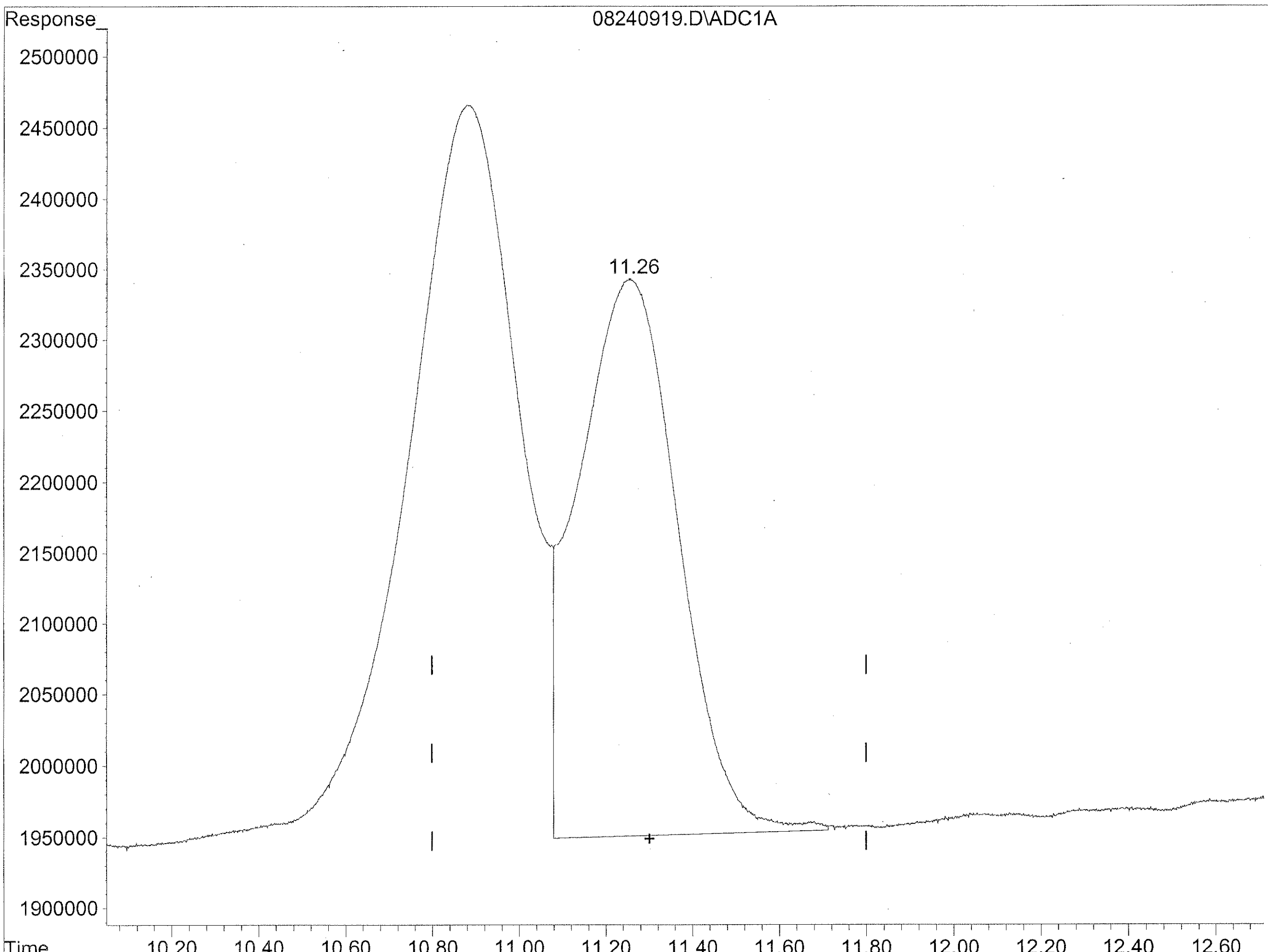
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.21	252393331	1374.831	ng/ml
2)	Acetaldehyde	1.70	190531159	1358.768	ng/ml
3)	Propionaldehyde	3.06	143656404	1346.418	ng/ml
4)	Crotonaldehyde	4.34	128230832	1316.333	ng/ml
5)	Butyraldehyde	5.24	120424860	1363.257	ng/ml
6)	Benzaldehyde	6.69	89082574	1352.414	ng/ml
7)	Isovaleraldehyde	7.54	104524706	1335.762	ng/ml
8)	Valeraldehyde	8.00	94875521	1290.735	ng/ml
9)	o-Tolualdehyde	8.83	79900971	1370.034	ng/ml
10)	m,p-Tolualdehyde	9.14	144989124	2685.213	ng/ml
11)	Hexaldehyde	10.88	92737947	1377.083	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.26	62609370	1277.393	ng/mlm

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240919.D   Vial: 17
Acq On : 24 Aug 2009 5:01 pm   Operator: HC
Sample : CCV 1500ng/ml S21-08240901   Inst : LC 01
Misc :   Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 11:34 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 12:41:27 2009
Response via : Multiple Level Calibration



QEdit

(12) 2,5-Dimethylbenzaldehyde

11.26min 1277.393ng/ml m

response 62609370

247

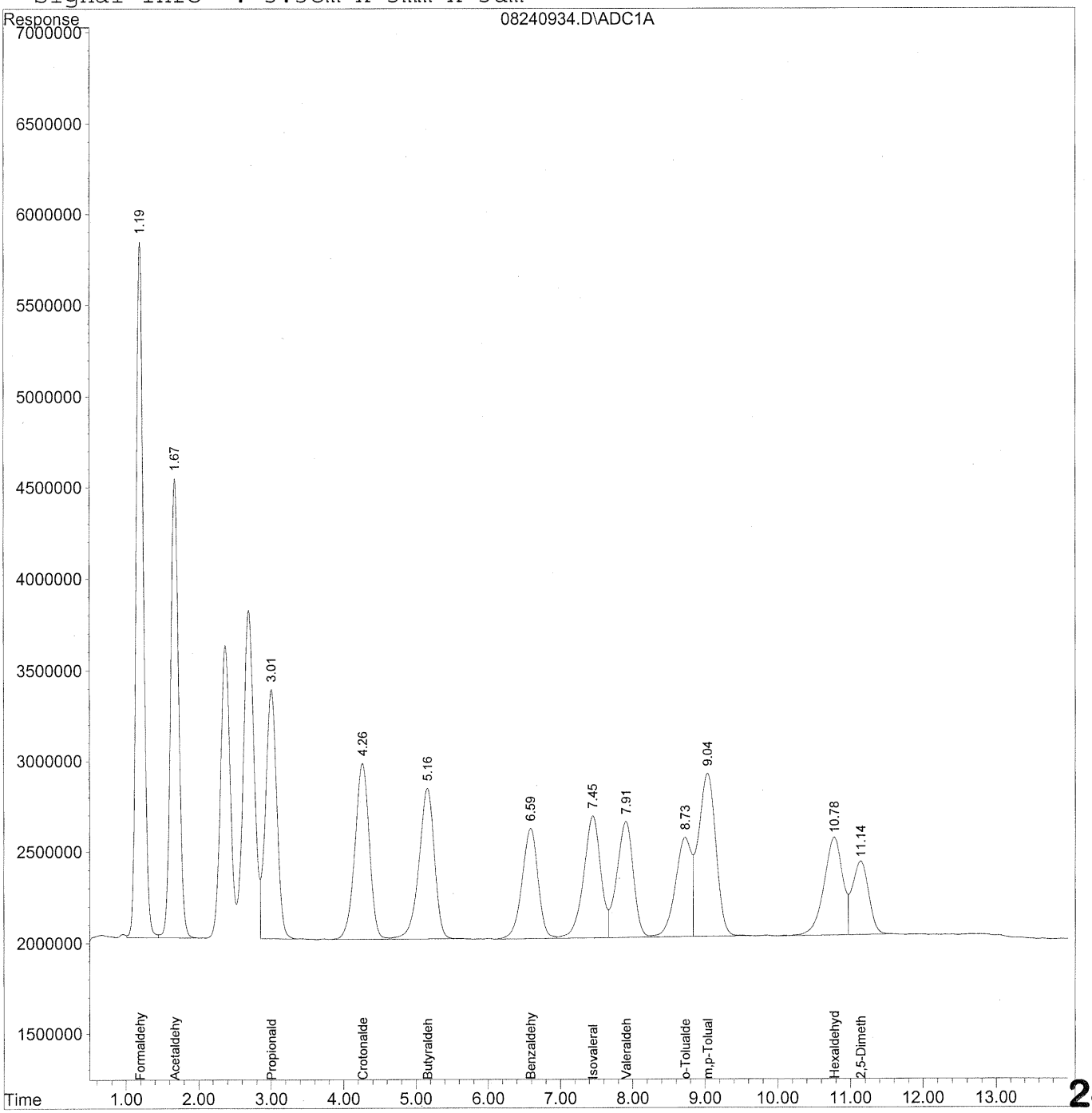
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240934.D Vial: 32
Acq On : 24 Aug 2009 8:47 pm Operator: HC
Sample : CCV 1500ng/ml S21-08240901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:23 19109 Quant Results File: TO110709.RES

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



248

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\24\08240934.D Vial: 32
Acq On : 24 Aug 2009 8:47 pm Operator: HC
Sample : CCV 1500ng/ml S21-08240901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:23 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 2009
Response via : Initial Calibration
DataAcq Meth : TO11S.M

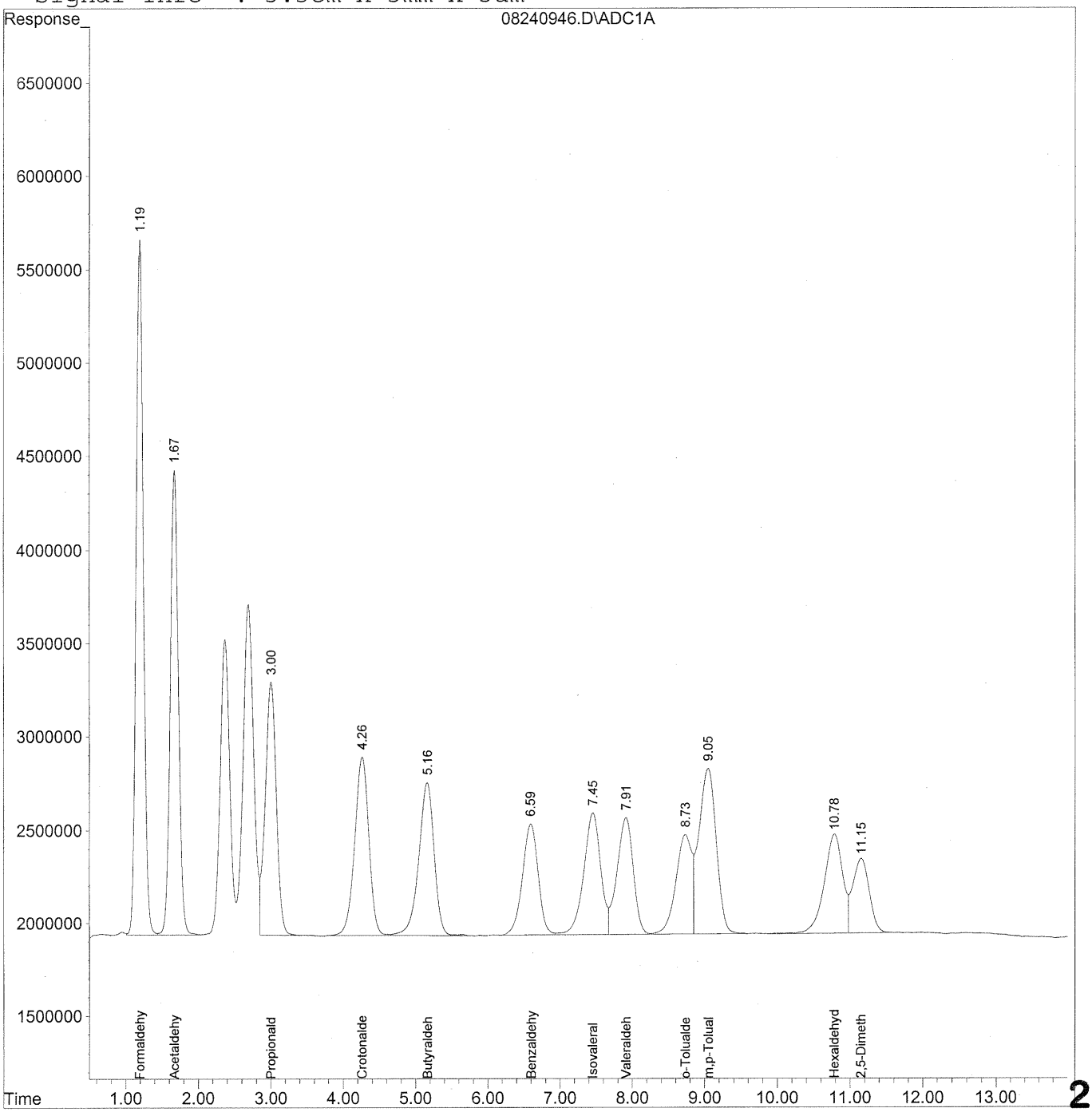
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.19	261168767	1422.632	ng/ml
2)	Acetaldehyde	1.67	197276586	1406.872	ng/ml
3)	Propionaldehyde	3.00	148406019	1390.934	ng/ml
4)	Crotonaldehyde	4.26	132550674	1360.678	ng/ml
5)	Butyraldehyde	5.16	124208971	1406.095	ng/ml
6)	Benzaldehyde	6.59	92307393	1401.372	ng/ml
7)	Isovaleraldehyde	7.45	110018810	1405.973	ng/ml
8)	Valeraldehyde	7.91	100336922	1365.035	ng/ml
9)	o-Tolualdehyde	8.72	83606258	1433.568	ng/ml
10)	m,p-Tolualdehyde	9.04	151429041	2804.480	ng/ml
11)	Hexaldehyde	10.78	95937308	1424.591	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.14	62516671	1275.502	ng/ml

Data File : J:\LC01\DATA\TO11\2009_08\24\08240946.D Vial: 43
Acq On : 24 Aug 2009 11:47 pm Operator: HC
Sample : CCV 1500ng/ml S21-08240901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:23 19109 Quant Results File: TO110709.RES

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



250

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\24\08240946.D Vial: 43
Acq On : 24 Aug 2009 11:47 pm Operator: HC
Sample : CCV 1500ng/ml S21-08240901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:23 19109 Quant Results File: TO110709.RES

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

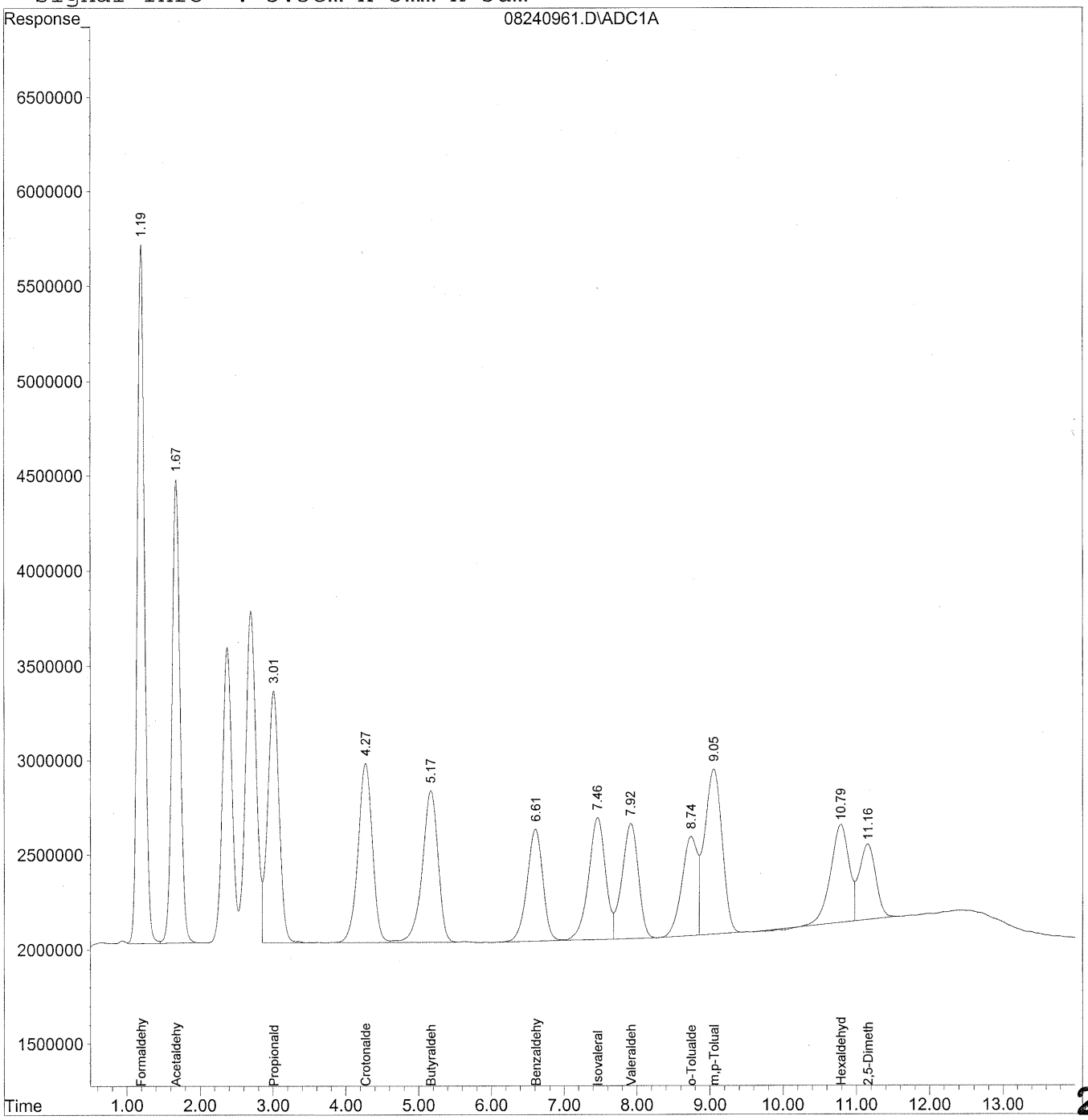
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.19	253995864	1383.560	ng/ml
2)	Acetaldehyde	1.67	193972280	1383.308	ng/ml
3)	Propionaldehyde	3.00	145980213	1368.198	ng/ml
4)	Crotonaldehyde	4.26	130950685	1344.253	ng/ml
5)	Butyraldehyde	5.16	123341530	1396.275	ng/ml
6)	Benzaldehyde	6.60	90261865	1370.317	ng/ml
7)	Isovaleraldehyde	7.45	107584035	1374.858	ng/ml
8)	Valeraldehyde	7.91	98568415	1340.975	ng/ml
9)	o-Tolualdehyde	8.73	82381082	1412.560	ng/ml
10)	m,p-Tolualdehyde	9.05	149515504	2769.042	ng/ml
11)	Hexaldehyde	10.78	94977473	1410.338	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.15	63586976	1297.339	ng/ml

Quantitation Report

Data File : J:\LC01\DATA\TO11\2009_08\24\08240961.D Vial: 58
Acq On : 25 Aug 2009 3:33 am Operator: HC
Sample : CCV 1500ng/ml S21-08240901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

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



252

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2009_08\24\08240961.D Vial: 58
Acq On : 25 Aug 2009 3:33 am Operator: HC
Sample : CCV 1500ng/ml S21-08240901 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Aug 25 7:20 19109 Quant Results File: TO110709.RES

Quant Method : J:\LC01\METHODS\TO110709.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Aug 24 08:44:34 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	1.19	250585282	1364.982	ng/ml
2)	Acetaldehyde	1.67	190792170	1360.629	ng/ml
3)	Propionaldehyde	3.01	145128630	1360.217	ng/ml
4)	Crotonaldehyde	4.27	128618213	1320.310	ng/ml
5)	Butyraldehyde	5.17	120433440	1363.354	ng/ml
6)	Benzaldehyde	6.61	88603389	1345.139	ng/ml
7)	Isovaleraldehyde	7.46	103869827	1327.393	ng/ml
8)	Valeraldehyde	7.92	94220063	1281.818	ng/ml
9)	o-Tolualdehyde	8.74	78227370	1341.338	ng/ml
10)	m,p-Tolualdehyde	9.05	145085056	2686.989	ng/ml
11)	Hexaldehyde	10.79	89794153	1333.370	ng/ml
12)	2,5-Dimethylbenzaldehyde	11.15	62668637	1278.602	ng/ml

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

255

Injection Log

Directory: j:\lc01\data\to11\2009_08\24

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	1	08240901.d	1.	PRIME		24 Aug 109 13::
2	2	08240902.d	1.	1500ng/ml TO11A std S21-08240901		24 Aug 109 13::
3	3	08240903.d	1.	ACN blank Lot CY023		24 Aug 109 12::
4	4	08240904.d	1.	P0902878-013 front 10x		24 Aug 109 12::
5	5	08240905.d	1.	P0902878-014 front 10x		24 Aug 109 12::
6	6	08240906.d	1.	P0902878-016 front 10x		24 Aug 109 12::
7	7	08240907.d	1.	P0902878-017 front 10x		24 Aug 109 12::
8	8	08240908.d	1.	CCV 1500ng/ml S21-08240901		24 Aug 109 12::
9	9	08240909.d	1.	CCV 1500ng/ml S21-08240901		24 Aug 109 12::
10	10	08240910.d	1.	ACN blk lot CY023		24 Aug 109 12::
11	10	08240911.d	1.	MB Front LOT 5899 1.0ml		24 Aug 109 12::
12	11	08240912.d	1.	MB Back LOT 5899 1.0ml		24 Aug 109 12::
13	12	08240913.d	1.	P0902901-001 back 1.0ml		24 Aug 109 12::
14	13	08240914.d	1.	P0902901-002 back 1.0ml		24 Aug 109 12::
15	14	08240915.d	1.	P0902901-001 front 1.0ml		24 Aug 109 12::
16	14	08240916.d	1.	P0902901-001dup front 1.0ml		24 Aug 109 12::
17	15	08240917.d	1.	P0902901-002 front 1.0ml		24 Aug 109 12::
18	16	08240918.d	1.	ACN wash		24 Aug 109 12::
19	17	08240919.d	1.	CCV 1500ng/ml S21-08240901		24 Aug 109 12::
20	18	08240920.d	1.	ACN lot Blk CY023		24 Aug 109 12::
21	19	08240921.d	1.	MB front lot 5855/5994 1.0ml		24 Aug 109 12::
22	20	08240922.d	1.	MB back lot 5855/5994 1.0ml		24 Aug 109 12::
23	21	08240923.d	1.	P0902912-001 back 1.0ml		24 Aug 109 12::
24	22	08240924.d	1.	P0902912-002 back 1.0ml		24 Aug 109 12::
25	23	08240925.d	1.	P0902912-003 back 1.0ml		24 Aug 109 12::
26	24	08240926.d	1.	P0902912-004 back 1.0ml		24 Aug 109 12::
27	25	08240927.d	1.	P0902912-005 back 1.0ml		24 Aug 109 12::
28	26	08240928.d	1.	P0902910-001 back 1.0ml		24 Aug 109 12::
29	27	08240929.d	1.	P0902910-002 back 1.0ml		24 Aug 109 12::
30	28	08240930.d	1.	P0902910-003 back 1.0ml		24 Aug 109 12::
31	29	08240931.d	1.	P0902910-004 back 1.0ml		24 Aug 109 12::
32	30	08240932.d	1.	P0902910-005 back 1.0ml		24 Aug 109 12::
33	31	08240933.d	1.	ACN wash		24 Aug 109 12::
34	32	08240934.d	1.	CCV 1500ng/ml S21-08240901		24 Aug 109 12::
35	33	08240935.d	1.	P0902910-006 back 1.0ml		24 Aug 109 12::
36	33	08240936.d	1.	P0902910-006dup back 1.0ml		24 Aug 109 12::
37	34	08240937.d	1.	P0902910-007 back 1.0ml		24 Aug 109 12::
38	35	08240938.d	1.	P0902910-008 back 1.0ml		24 Aug 109 12::
39	36	08240939.d	1.	P0902910-009 back 1.0ml		24 Aug 109 13::
40	37	08240940.d	1.	P0902910-010 back 1.0ml		24 Aug 109 13::
41	38	08240941.d	1.	P0902910-011 back 1.0ml		24 Aug 109 13::
42	39	08240942.d	1.	P0902910-012 back 1.0ml		24 Aug 109 13::
43	40	08240943.d	1.	P0902910-013 back 1.0ml		24 Aug 109 13::
44	41	08240944.d	1.	P0902910-014 back 1.0ml		24 Aug 109 13::
45	42	08240945.d	1.	ACN wash		24 Aug 109 13::
46	43	08240946.d	1.	CCV 1500ng/ml S21-08240901		24 Aug 109 13::
47	44	08240947.d	1.	ACN blk lot CY023		25 Aug 109 13::
48	45	08240948.d	1.	MB front lot 5855/5994 1.0ml		25 Aug 109 13::
49	46	08240949.d	1.	MB back lot 5855/5994 1.0ml		25 Aug 109 13::
50	47	08240950.d	1.	P0902910-015 back 1.0ml		25 Aug 109 13::
51	48	08240951.d	1.	P0902910-016 back 1.0ml		25 Aug 109 12::
52	49	08240952.d	1.	P0902910-017 back 1.0ml		25 Aug 109 12::
53	50	08240953.d	1.	P0902910-018 back 1.0ml		25 Aug 109 12::
54	51	08240954.d	1.	P0902912-001 front 1.0ml		25 Aug 109 12::
55	52	08240955.d	1.	P0902912-002 front 1.0ml		25 Aug 109 12::
56	53	08240956.d	1.	P0902912-003 front 1.0ml		25 Aug 109 12::
57	54	08240957.d	1.	P0902912-004 front 1.0ml		25 Aug 109 12::

Injection Log

Directory: j:\lc01\data\to11\2009_08\24

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
58	55	08240958.d	1.	P0902912-005 front 1.0ml		25 Aug 109 12::
59	56	08240959.d	1.	P0902910-001 front 1.0ml		25 Aug 109 12::
60	57	08240960.d	1.	ACN wash		25 Aug 109 12::
61	58	08240961.d	1.	CCV 1500ng/ml S21-08240901		25 Aug 109 12::
62	59	08240962.d	1.	P0902910-002 front 1.0ml		25 Aug 109 12::
63	59	08240963.d	1.	P0902910-002dup front 1.0ml		25 Aug 109 12::
64	60	08240964.d	1.	P0902910-003 front 1.0ml		25 Aug 109 12::
65	61	08240965.d	1.	P0902910-004 front 1.0ml		25 Aug 109 12::
66	62	08240966.d	1.	P0902910-005 front 1.0ml		25 Aug 109 12::
67	63	08240967.d	1.	P0902910-006 front 1.0ml		25 Aug 109 12::
68	64	08240968.d	1.	P0902910-007 front 1.0ml		25 Aug 109 12::
69	65	08240969.d	1.	P0902910-008 front 1.0ml		25 Aug 109 12::
70	66	08240970.d	1.	P0902910-009 front 1.0ml		25 Aug 109 12::
71	67	08240971.d	1.	P0902910-010 front 1.0ml		25 Aug 109 12::
72	68	08240972.d	1.	ACN wash		25 Aug 109 12::
73	69	08240973.d	1.	CCV 1500ng/ml S21-08240901		25 Aug 109 12::
74	70	08240974.d	1.	ACN blk lot CY023		25 Aug 109 12::
75	71	08240975.d	1.	MB front lot 5855/5994 1.0ml		25 Aug 109 12::
76	72	08240976.d	1.	MB back lot 5855/5994 1.0ml		25 Aug 109 12::
77	73	08240977.d	1.	P0902910-011 front 1.0ml		25 Aug 109 12::
78	73	08240978.d	1.	P0902910-011dup front 1.0ml		25 Aug 109 12::
79	74	08240979.d	1.	P0902910-012 front 1.0ml		25 Aug 109 12::
80	75	08240980.d	1.	P0902910-013 front 1.0ml		25 Aug 109 12::
81	76	08240981.d	1.	P0902910-014 front 1.0ml		25 Aug 109 12::
82	77	08240982.d	1.	P0902910-015 front 1.0ml		25 Aug 109 12::
83	78	08240983.d	1.	P0902910-016 front 1.0ml		25 Aug 109 12::
84	79	08240984.d	1.	P0902910-017 front 1.0ml		25 Aug 109 12::
85	80	08240985.d	1.	P0902910-018 front 1.0ml		25 Aug 109 12::
86	81	08240986.d	1.	ACN wash		25 Aug 109 12::
87	82	08240987.d	1.	CCV 1500ng/ml S21-08240901		25 Aug 109 13::
88	83	08240988.d	1.	ACN CY023		25 Aug 109 13::
89	84	08240989.d	1.	P0902910-015 front 10x		25 Aug 109 13::
90	85	08240990.d	1.	CCV 1500ng/ml S21-08240901		25 Aug 109 13::