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

July 16, 2010

Data Coordinator
Environmental Health & Engineering, Inc.
117 Fourth Avenue
Needham, MA 02494

RE: 17131

Dear Data Coordinator:

Enclosed are the results of the samples submitted to our laboratory on June 19, 2010. For your reference, these analyses have been assigned our service request number P1002127.

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 225 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; United States Department of Defense Environmental Laboratory Accreditation Program (DoD-ELAP), Certificate No. L10-3; Pennsylvania Registration No. 68-03307; TX Commission of Environmental Quality, NELAP ID T104704413-09-TX; Minnesota Department of Health, Certificate No. 11495AA; Washington State Department of Ecology, ELAP Lab ID: C946. 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
Project Manager

Client: Environmental Health & Engineering, Inc. CAS Project No: P1002127
Project: 17131

CASE NARRATIVE

The samples were received intact under chain of custody on June 19, 2010 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.

Formaldehyde Analysis

The samples were analyzed for formaldehyde 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, Inc.
Project: 17131

Service Request: P1002127

SAMPLE CROSS-REFERENCE

SAMPLE #	CLIENT SAMPLE ID	DATE	TIME
P1002127-001	111383	6/18/10	00:00
P1002127-002	111384	6/18/10	00:00
P1002127-003	111385	6/18/10	00:00
P1002127-004	111386	6/18/10	00:00
P1002127-005	111387	6/18/10	00:00
P1002127-006	111399	6/18/10	00:00
P1002127-007	111400	6/18/10	00:00
P1002127-008	111401	6/18/10	00:00
P1002127-009	111402	6/18/10	00:00
P1002127-010	111403	6/18/10	00:00
P1002127-011	111302	6/18/10	00:00
P1002127-012	111303	6/18/10	00:00
P1002127-013	111304	6/18/10	00:00
P1002127-014	111305	6/18/10	00:00
P1002127-015	111306	6/18/10	00:00
P1002127-016	111317	6/18/10	00:00
P1002127-017	111318	6/18/10	00:00
P1002127-018	111319	6/18/10	00:00
P1002127-019	111320	6/18/10	00:00
P1002127-020	111321	6/18/10	00:00
P1002127-021	111332	6/18/10	00:00
P1002127-022	111333	6/18/10	00:00
P1002127-023	111334	6/18/10	00:00
P1002127-024	111335	6/18/10	00:00
P1002127-025	111336	6/18/10	00:00
P1002127-026	111347	6/18/10	00:00
P1002127-027	111348	6/18/10	00:00
P1002127-028	111349	6/18/10	00:00
P1002127-029	111350	6/18/10	00:00
P1002127-030	111351	6/18/10	00:00

Environmental Health & Engineering, Inc.

CHAIN OF CUSTODY FORM

DATE: 18 JUN 10

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

P1002127

TO: COLUMBIA

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 # 17131

The cost of this analysis will be covered by EH&E Purchase Order # 17131

For EH & E Data Coordinator - URGENT DATA ☐

	SAMPLE ID	SAMPLE TYPE	ANALYTICAL METHOD/NUMBER	OTHER (Time/Date/Vol.
1	111383	AIR	FORMALDEHYDE ANALYSIS ONLY	13D 23H 25M
2	111384	↓	↓	↓
3	111385			
4	111386			
5	111387			∅
6	111399			14D 15M
7	111400			↓
8	111401			
9	111402			
10	111403			∅
11	111302			13 DAYS
12	111303			↓
13	111304			
14	111305			
15	111306			∅
16	111317	↓	↓	13D 23H 50M

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 bbaker@eheinc.com

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

Relinquished by: _____ of Environmental Health & Engineering, Inc. Date: 6/18/10

Received by: _____ of (company name) CAT Date: 6/18/10 1050

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

Environmental Health & Engineering, Inc.

CHAIN OF CUSTODY FORM

DATE: 18 JUN 10

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

P1002127

TO: COLUMBIA

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 # 17131

The cost of this analysis will be covered by EH&E Purchase Order # 17131

For EH & E Data Coordinator - URGENT DATA ☒

	SAMPLE ID	SAMPLE TYPE	ANALYTICAL METHOD/NUMBER	OTHER: Time/Date/Vol.
17	111318	AIR	FORMALDEHYDE ANALYSIS ONLY	13D 23H 50M
18	111319			"
19	111320			"
20	111321			Ø
21	111332			13D 23H 10M
22	111333			"
23	111334			"
24	111335			"
25	111336			Ø
26	111347			13D 23H 40M
27	111348			"
28	111349			"
29	111350			"
30	111351			Ø

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 bbaker@eheinc.com

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

Relinquished by: _____ of Environmental Health & Engineering, Inc. Date: 6/18/10

Received by: _____ of (company name) CLS Date: 6/18/10 1050

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

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

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

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

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

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

Columbia Analytical Services, Inc.
Sample Acceptance Check Form

Client: Environmental Health & Engineering, Inc. Work order: P1002127

Project: 17131

Sample(s) received on: 06/19/10 Date opened: 06/19/10 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 2 °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
P1002127-001.01	Passive (Radiello DNPH)					
P1002127-002.01	Passive (Radiello DNPH)					
P1002127-003.01	Passive (Radiello DNPH)					
P1002127-004.01	Passive (Radiello DNPH)					
P1002127-005.01	Passive (Radiello DNPH)					

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

Chain of Custody is missing time collected

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

6

Columbia Analytical Services, Inc.
Sample Acceptance Check Form

Client: Environmental Health & Engineering, Inc. Work order: P1002127

Project: 17131

Sample(s) received on: 06/19/10 Date opened: 06/19/10 by: MZAMORA

Lab Sample ID	Container Description	Required pH *	Received pH	Adjusted pH	VOA Headspace (Presence/Absence)	Receipt / Preservation Comments
P1002127-006.01	Passive (Radiello DNPH)					
P1002127-007.01	Passive (Radiello DNPH)					
P1002127-008.01	Passive (Radiello DNPH)					
P1002127-009.01	Passive (Radiello DNPH)					
P1002127-010.01	Passive (Radiello DNPH)					
P1002127-011.01	Passive (Radiello DNPH)					
P1002127-012.01	Passive (Radiello DNPH)					
P1002127-013.01	Passive (Radiello DNPH)					
P1002127-014.01	Passive (Radiello DNPH)					
P1002127-015.01	Passive (Radiello DNPH)					
P1002127-016.01	Passive (Radiello DNPH)					
P1002127-017.01	Passive (Radiello DNPH)					
P1002127-018.01	Passive (Radiello DNPH)					
P1002127-019.01	Passive (Radiello DNPH)					
P1002127-020.01	Passive (Radiello DNPH)					
P1002127-021.01	Passive (Radiello DNPH)					
P1002127-022.01	Passive (Radiello DNPH)					
P1002127-023.01	Passive (Radiello DNPH)					
P1002127-024.01	Passive (Radiello DNPH)					
P1002127-025.01	Passive (Radiello DNPH)					
P1002127-026.01	Passive (Radiello DNPH)					
P1002127-027.01	Passive (Radiello DNPH)					
P1002127-028.01	Passive (Radiello DNPH)					
P1002127-029.01	Passive (Radiello DNPH)					
P1002127-030.01	Passive (Radiello DNPH)					

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

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

7

RESULTS OF ANALYSIS

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Client: **Environmental Health & Engineering, Inc.**
Client Project ID: **17131** CAS Project ID: P1002127

Formaldehyde

Test Code: EPA TO-11A
Instrument ID: HP1050/UV_Vis 360/LC2
Analyst: Madeleine Dangazyan
Sampling Media: Radiello Tube(s)
Test Notes: **BC**

Date(s) Collected: 6/18/10
Date Received: 6/19/10
Date Analyzed: 6/28 - 6/29/10
Desorption Volume: 2.0 ml(s)

Client Sample ID	CAS Sample ID	Sampling Time Minutes	Dilution Factor	**Result µg/Sample**	**Result µg/m³**	MRL µg/m³	**Result ppbV**	MRL ppbV	Data Qualifier
111383	P1002127-001	20125	10.0	**45**	**22**	1.0	**18**	0.82	
111384	P1002127-002	20125	10.0	**61**	**31**	1.0	**25**	0.82	
111385	P1002127-003	20125	10.0	**42**	**21**	1.0	**17**	0.82	
111386	P1002127-004	20125	1.0	**3.7**	**1.9**	0.10	**1.5**	0.082	
111387	P1002127-005	NA	1.0	**0.33**	NA	NA	NA	NA	
111399	P1002127-006	20175	10.0	**27**	**14**	1.0	**11**	0.82	
111400	P1002127-007	20175	10.0	**27**	**14**	1.0	**11**	0.82	
111401	P1002127-008	20175	10.0	**26**	**13**	1.0	**11**	0.82	
111402	P1002127-009	20175	1.0	**3.5**	**1.8**	0.10	**1.4**	0.082	
111403	P1002127-010	NA	1.0	**0.40**	NA	NA	NA	NA	
111302	P1002127-011	18720	10.0	**45**	**24**	1.1	**20**	0.88	
111303	P1002127-012	18720	10.0	**49**	**26**	1.1	**21**	0.88	
111304	P1002127-013	18720	10.0	**45**	**24**	1.1	**20**	0.88	
111305	P1002127-014	18720	1.0	**3.3**	**1.8**	0.11	**1.4**	0.088	
111306	P1002127-015	NA	1.0	**0.41**	NA	NA	NA	NA	
111317	P1002127-016	20150	10.0	**43**	**22**	1.0	**18**	0.82	

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
NA = Not applicable.
BC = Results reported are not blank corrected.

Verified By: Re Date: 7/12/10

Client: **Environmental Health & Engineering, Inc.**
Client Project ID: **17131** CAS Project ID: P1002127

Formaldehyde

Test Code: EPA TO-11A
Instrument ID: HP1050/UV_Vis 360/LC2
Analyst: Madeleine Dangazyan
Sampling Media: Radiello Tube(s)
Test Notes: **BC**

Date(s) Collected: 6/18/10
Date Received: 6/19/10
Date Analyzed: 6/28 - 6/29/10
Desorption Volume: 2.0 ml(s)

Client Sample ID	CAS Sample ID	Sampling Time Minutes	Dilution Factor	**Result µg/Sample**	**Result µg/m³**	MRL µg/m³	**Result ppbV**	MRL ppbV	Data Qualifier
111318	P1002127-017	20150	10.0	**48**	**24**	1.0	**20**	0.82	
111319	P1002127-018	20150	10.0	**47**	**23**	1.0	**19**	0.82	
111320	P1002127-019	20150	1.0	**3.7**	**1.8**	0.10	**1.5**	0.082	
111321	P1002127-020	NA	1.0	**0.39**	NA	NA	NA	NA	
111332	P1002127-021	20110	10.0	**43**	**21**	1.0	**17**	0.82	
111333	P1002127-022	20110	10.0	**43**	**22**	1.0	**18**	0.82	
111334	P1002127-023	20110	10.0	**42**	**21**	1.0	**17**	0.82	
111335	P1002127-024	20110	1.0	**3.2**	**1.6**	0.10	**1.3**	0.082	
111336	P1002127-025	NA	1.0	**0.40**	NA	NA	NA	NA	
111347	P1002127-026	20140	10.0	**46**	**23**	1.0	**19**	0.82	
111348	P1002127-027	20140	10.0	**51**	**26**	1.0	**21**	0.82	
111349	P1002127-028	20140	10.0	**67**	**34**	1.0	**28**	0.82	
111350	P1002127-029	20140	1.0	**3.7**	**1.9**	0.10	**1.5**	0.082	
111351	P1002127-030	NA	1.0	**0.42**	NA	NA	NA	NA	
Method Blank	P100628-MB	NA	1.0	< 0.20	NA	NA	NA	NA	

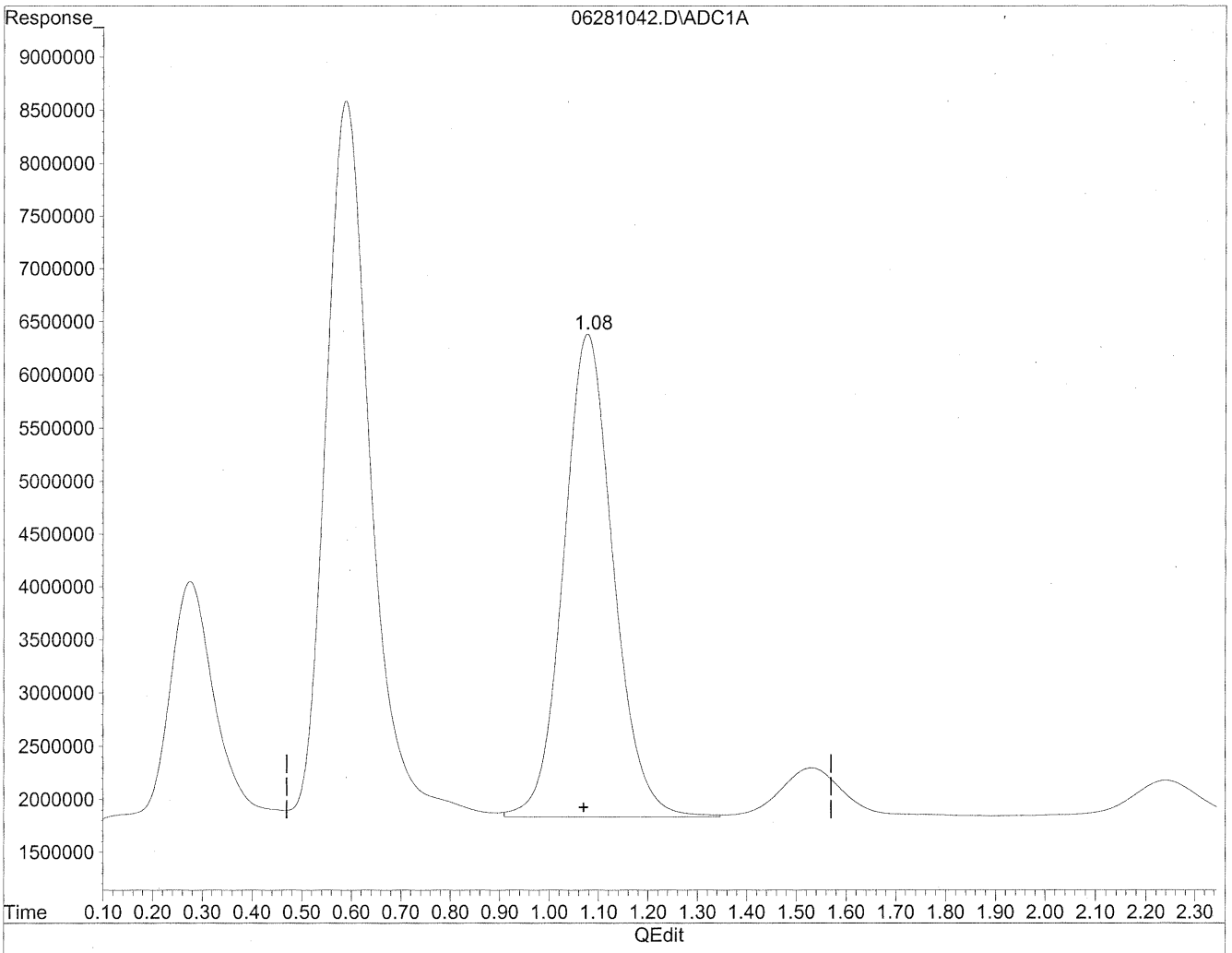
ND = Compound was analyzed for, but not detected above the laboratory reporting limit.
NA = Not applicable.
BC = Results reported are not blank corrected.

Verified By: Re Date: 7/12/10

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281042.D Vial: 42
Acq On : 28 Jun 2010 19:35 Operator: MD
Sample : P1002127-001 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:27 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 2225.081ng/ml

response 318599648

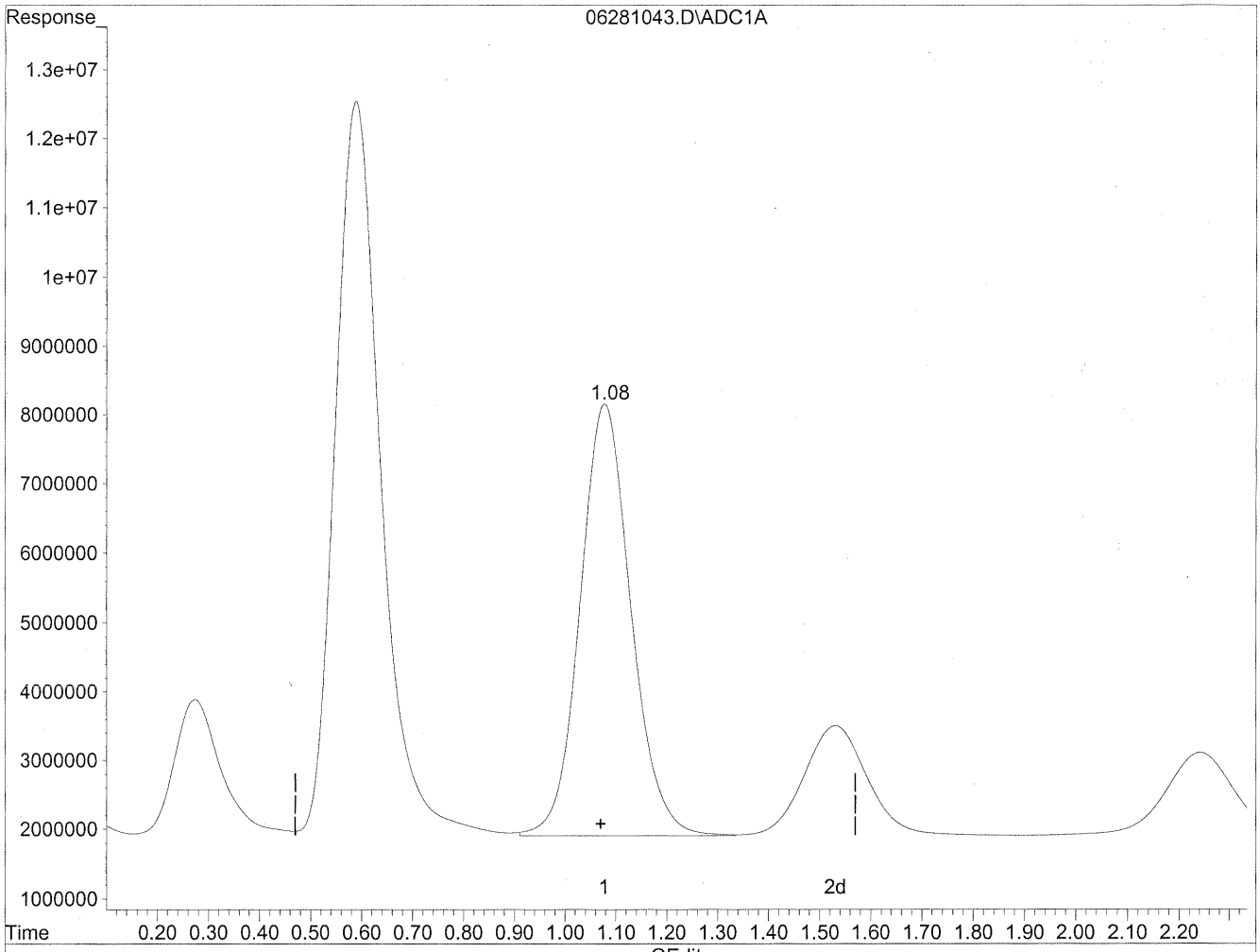
(+) = Expected Retention Time

06281042.D TO110610.M Tue Jun 29 13:27:59 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281043.D Vial: 43
Acq On : 28 Jun 2010 19:49 Operator: MD
Sample : P1002127-002 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:28 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 3040.650ng/ml

response 435377511

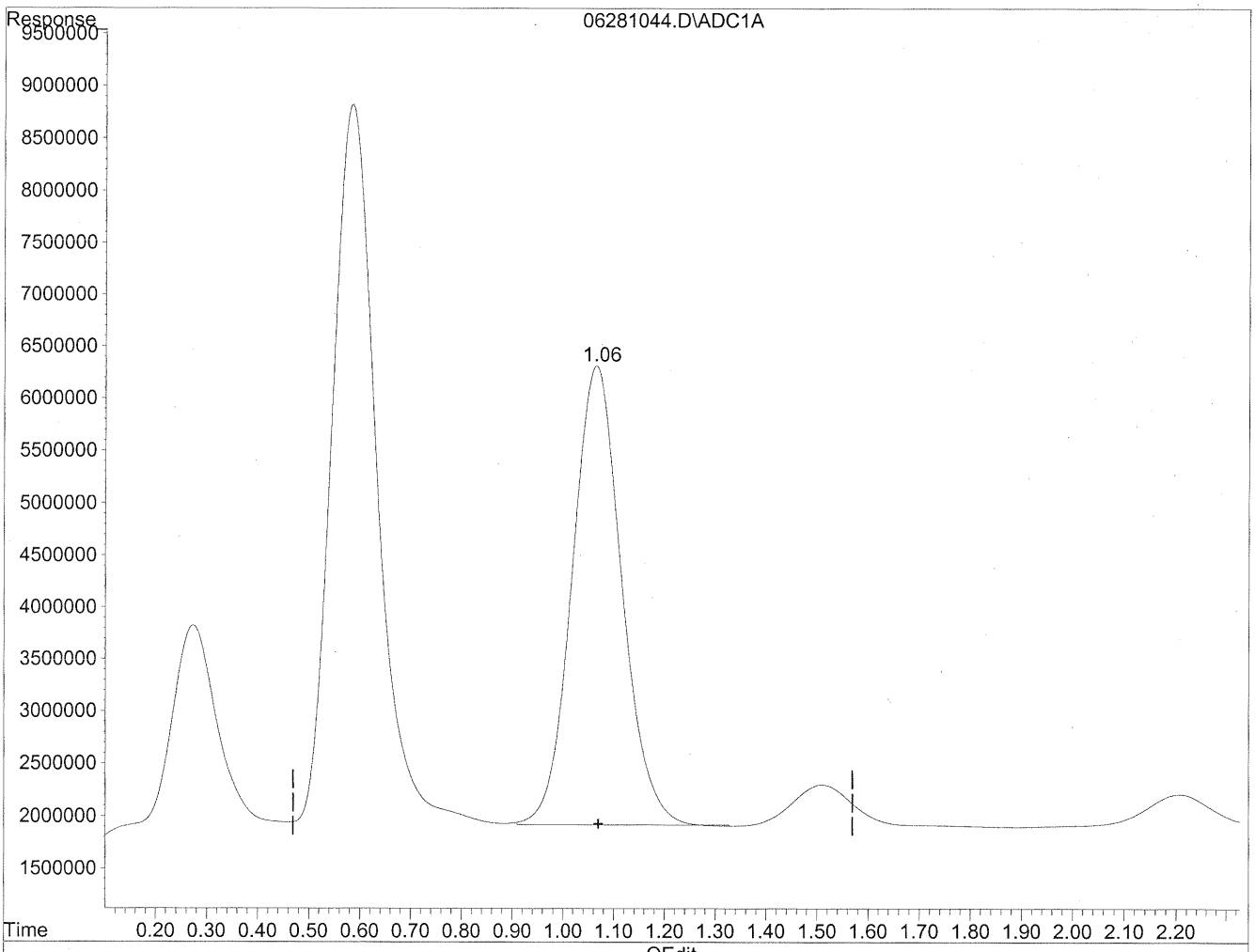
(+) = Expected Retention Time

06281043.D TO110610.M Tue Jun 29 13:28:09 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281044.D Vial: 44
Acq On : 28 Jun 2010 20:03 Operator: MD
Sample : P1002127-003 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:28 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.06min 2079.912ng/ml

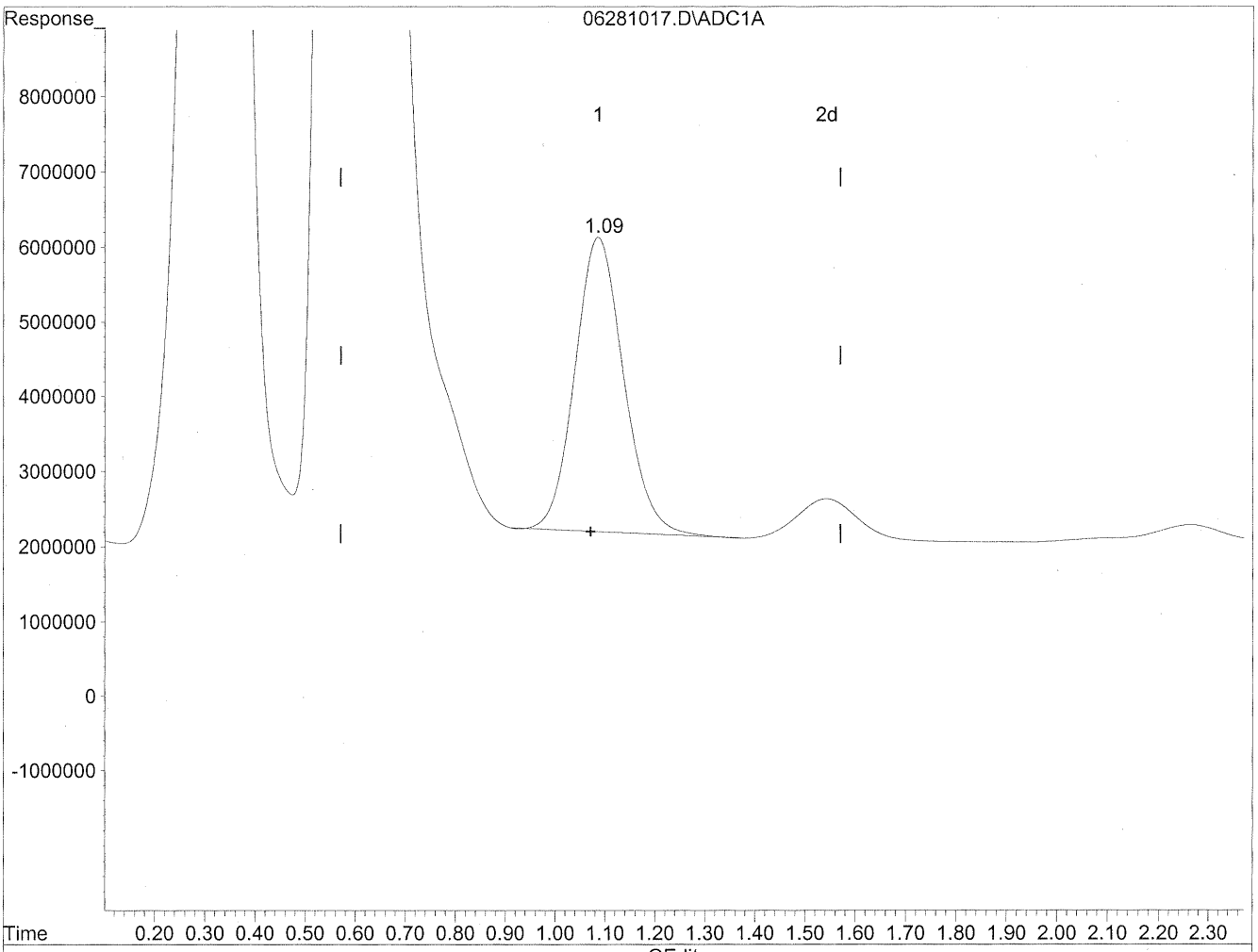
response 297813594

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281017.D Vial: 17
Acq On : 28 Jun 2010 13:57 Operator: MD
Sample : P1002127-004 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 28 14:18 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.09min 1866.219ng/ml

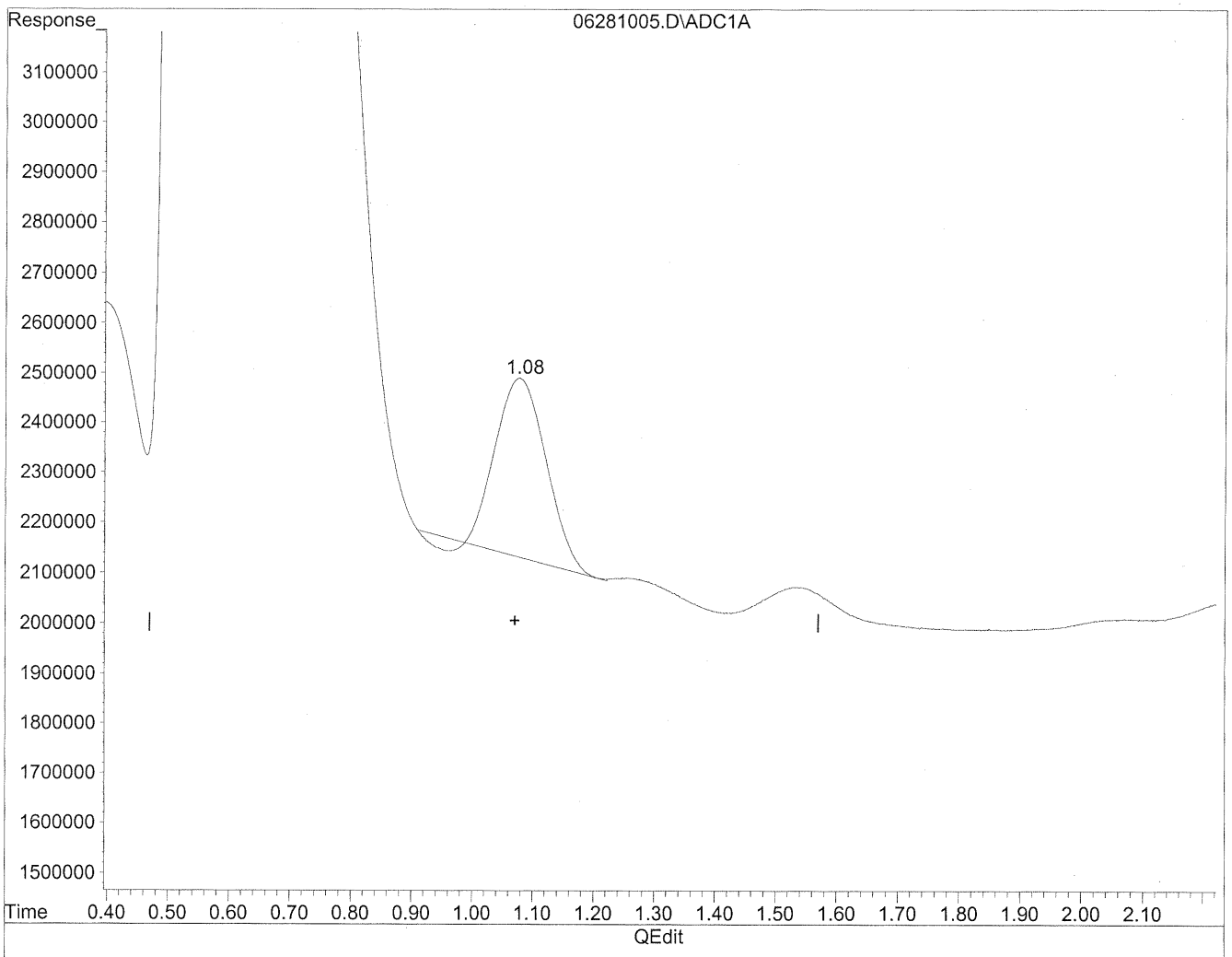
response 267215777

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281005.D Vial: 5
Acq On : 28 Jun 2010 11:01 Operator: MD
Sample : P1002127-005 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:18 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 140.865ng/ml

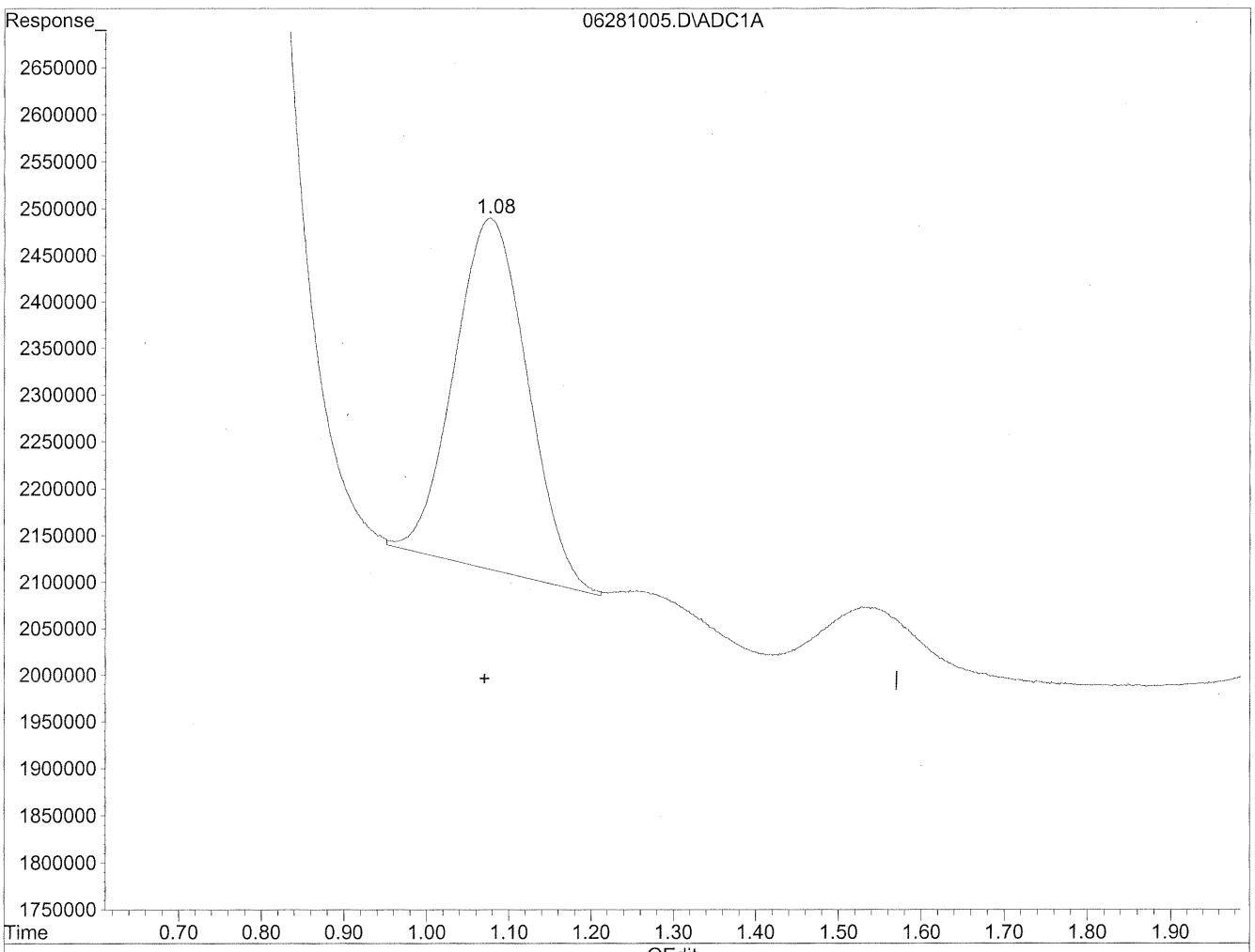
response 20169814

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281005.D Vial: 5
Acq On : 28 Jun 2010 11:01 Operator: MD
Sample : P1002127-005 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:18 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 162.696ng/ml m

response 23295796

6/30/10

6/30/10

16

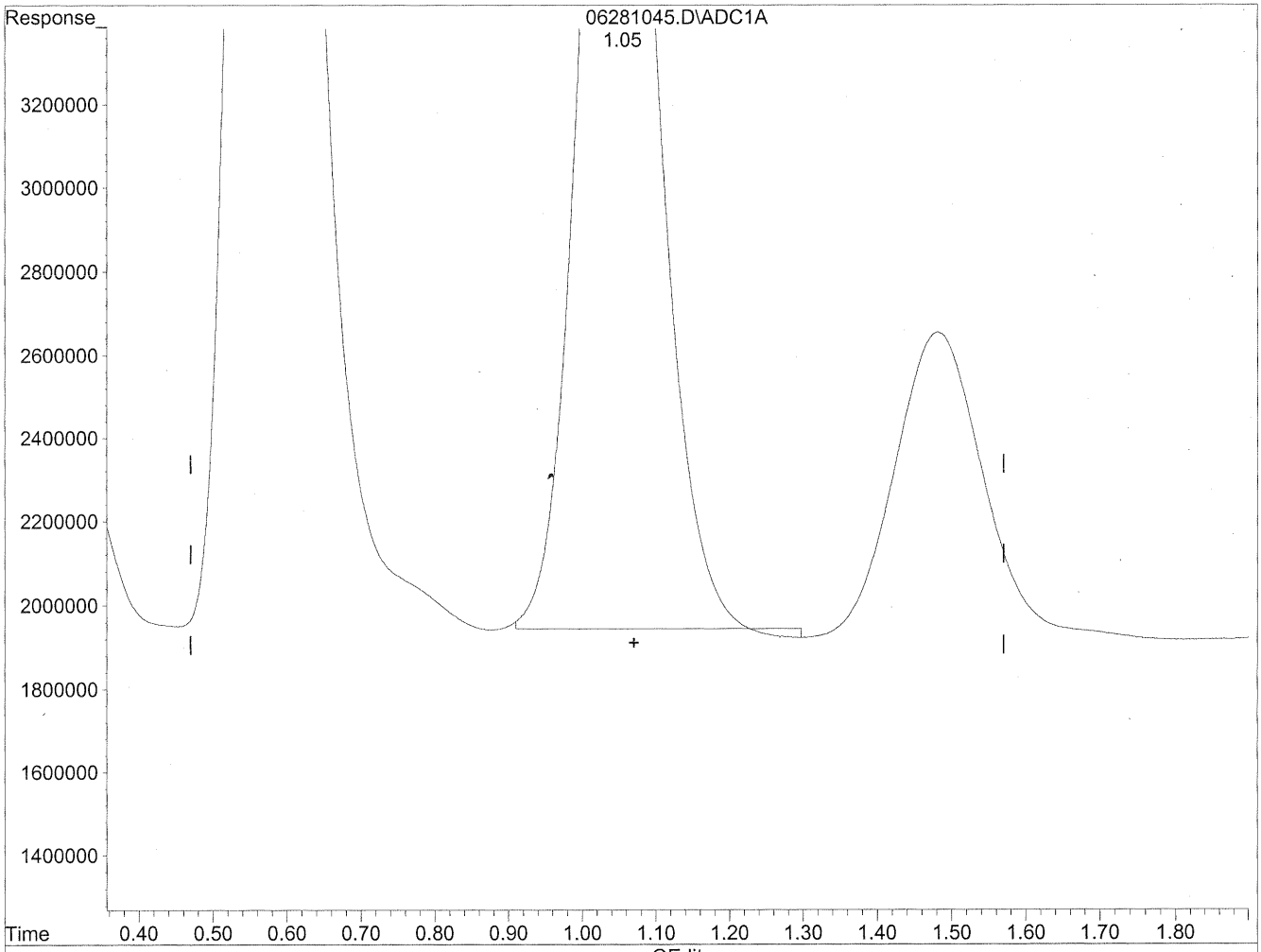
(+) = Expected Retention Time

06281005.D TO110610.M Tue Jun 29 13:20:18 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281045.D Vial: 45
Acq On : 28 Jun 2010 20:16 Operator: MD
Sample : P1002127-006 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:28 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.05min 1341.592ng/ml

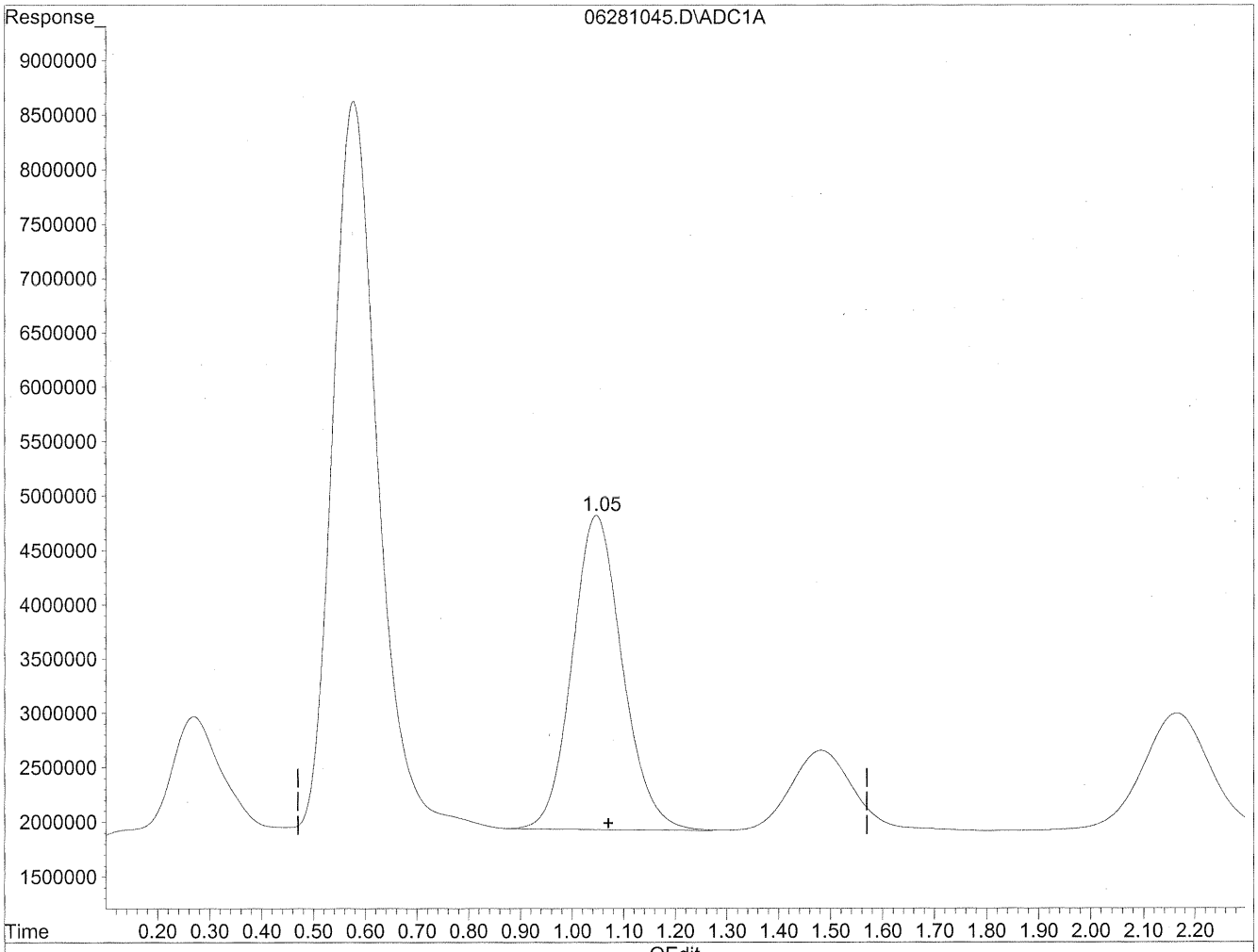
response 192096768

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281045.D Vial: 45
Acq On : 28 Jun 2010 20:16 Operator: MD
Sample : P1002127-006 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:28 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.05min 1366.450ng/ml m

response 195656023

6/30/10

6/30/10

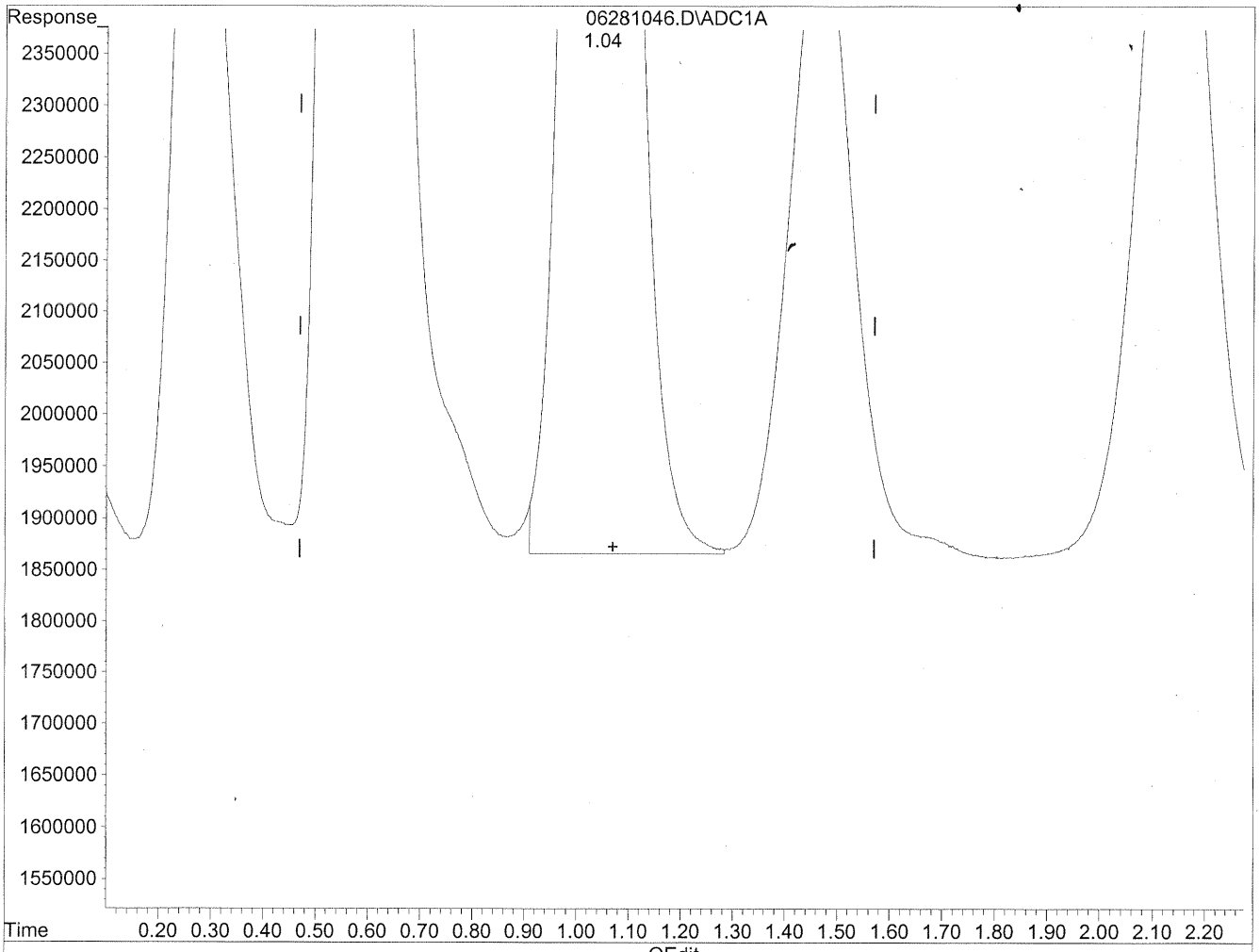
(+) = Expected Retention Time

06281045.D TO110610.M Tue Jun 29 13:28:36 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281046.D Vial: 46
Acq On : 28 Jun 2010 20:30 Operator: MD
Sample : P1002127-007 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:28 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.04min 1362.318ng/ml

response 195064350

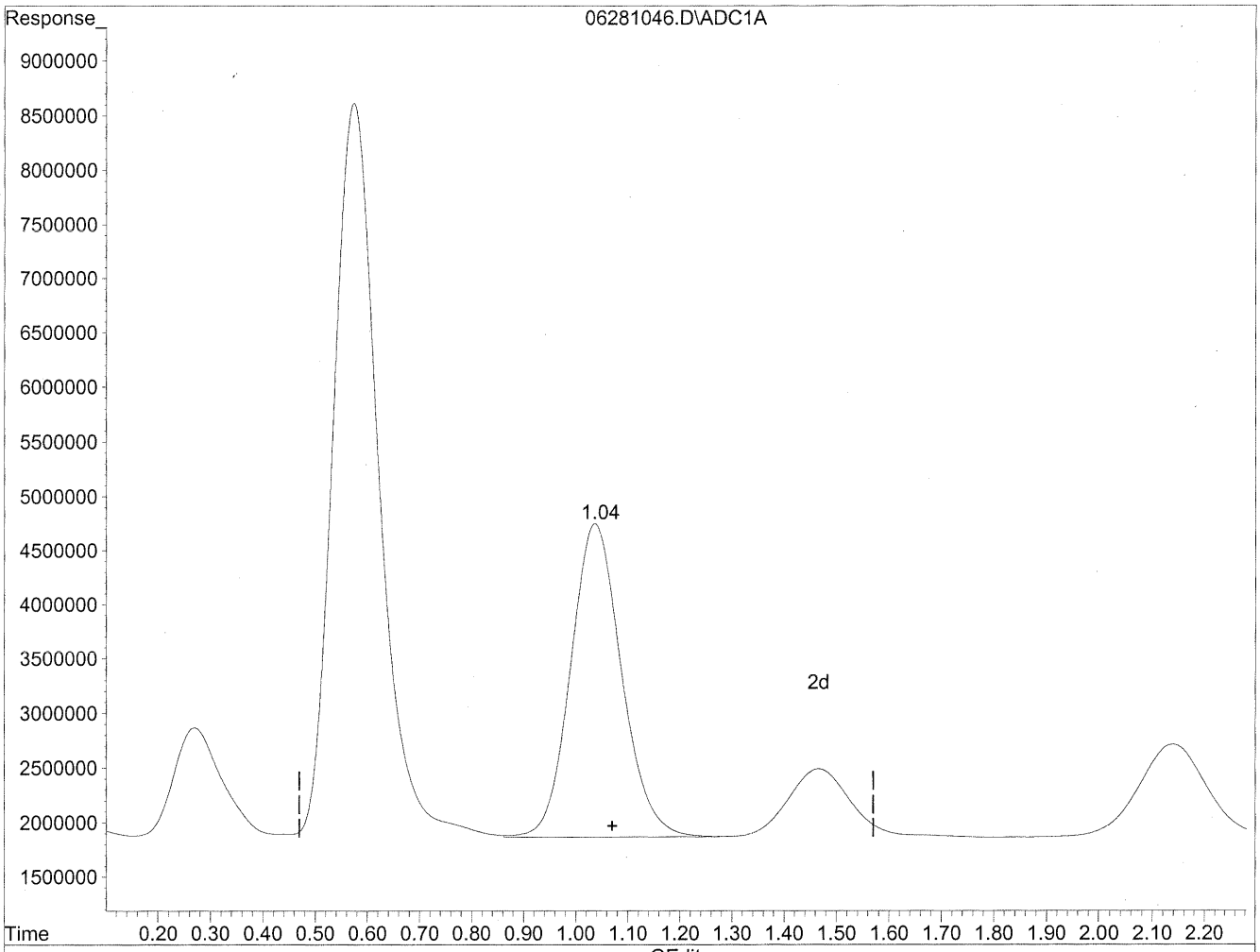
(+) = Expected Retention Time

06281046.D TO110610.M Tue Jun 29 14:56:07 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281046.D Vial: 46
Acq On : 28 Jun 2010 20:30 Operator: MD
Sample : P1002127-007 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:28 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.04min 1366.009ng/ml m

response 195592843

HC 6/29/10

IC 6/30/10

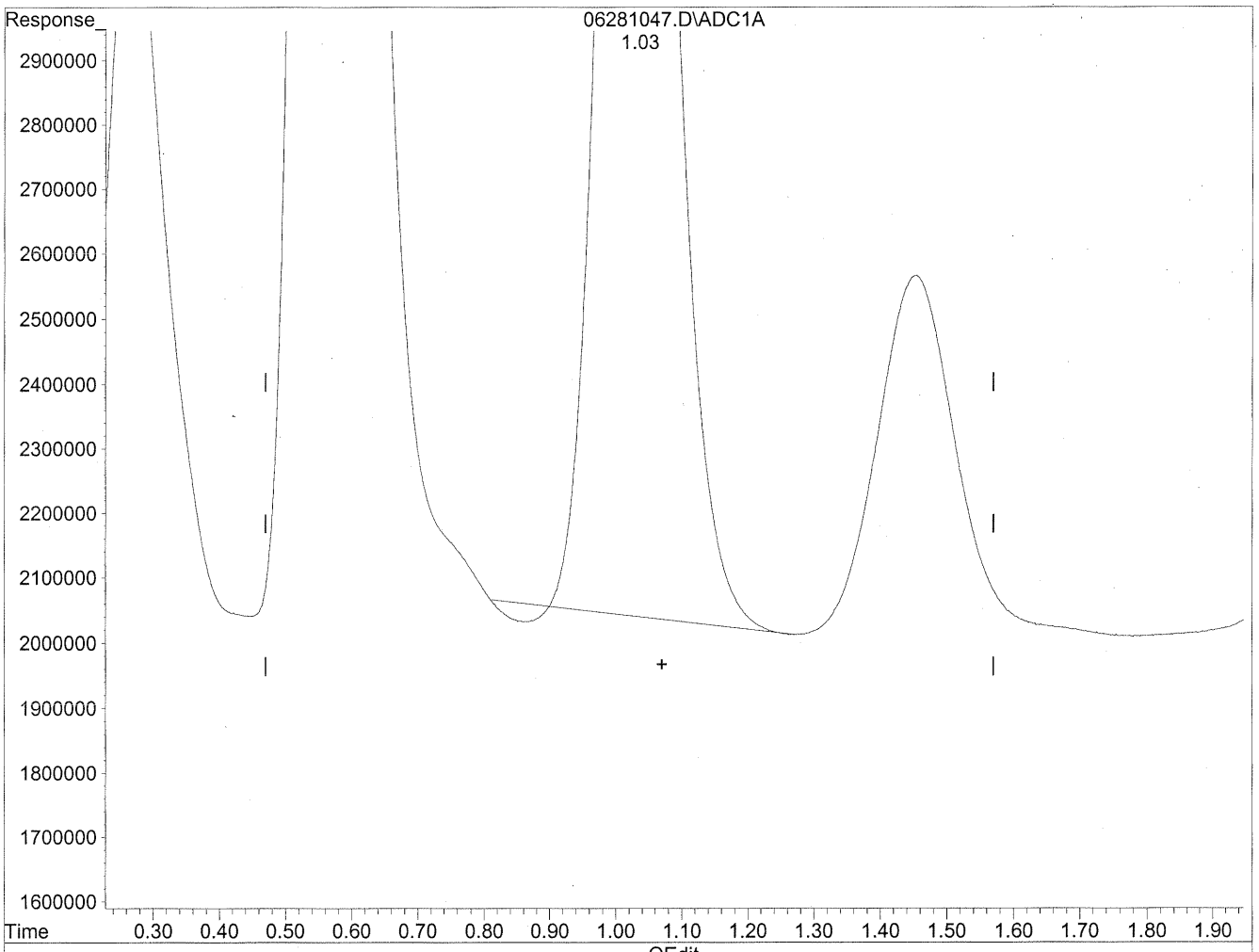
(+) = Expected Retention Time

06281046.D TO110610.M Tue Jun 29 14:56:13 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281047.D Vial: 47
Acq On : 28 Jun 2010 20:43 Operator: MD
Sample : P1002127-008 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 14:56 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.03min 1270.221ng/ml

response 181877383

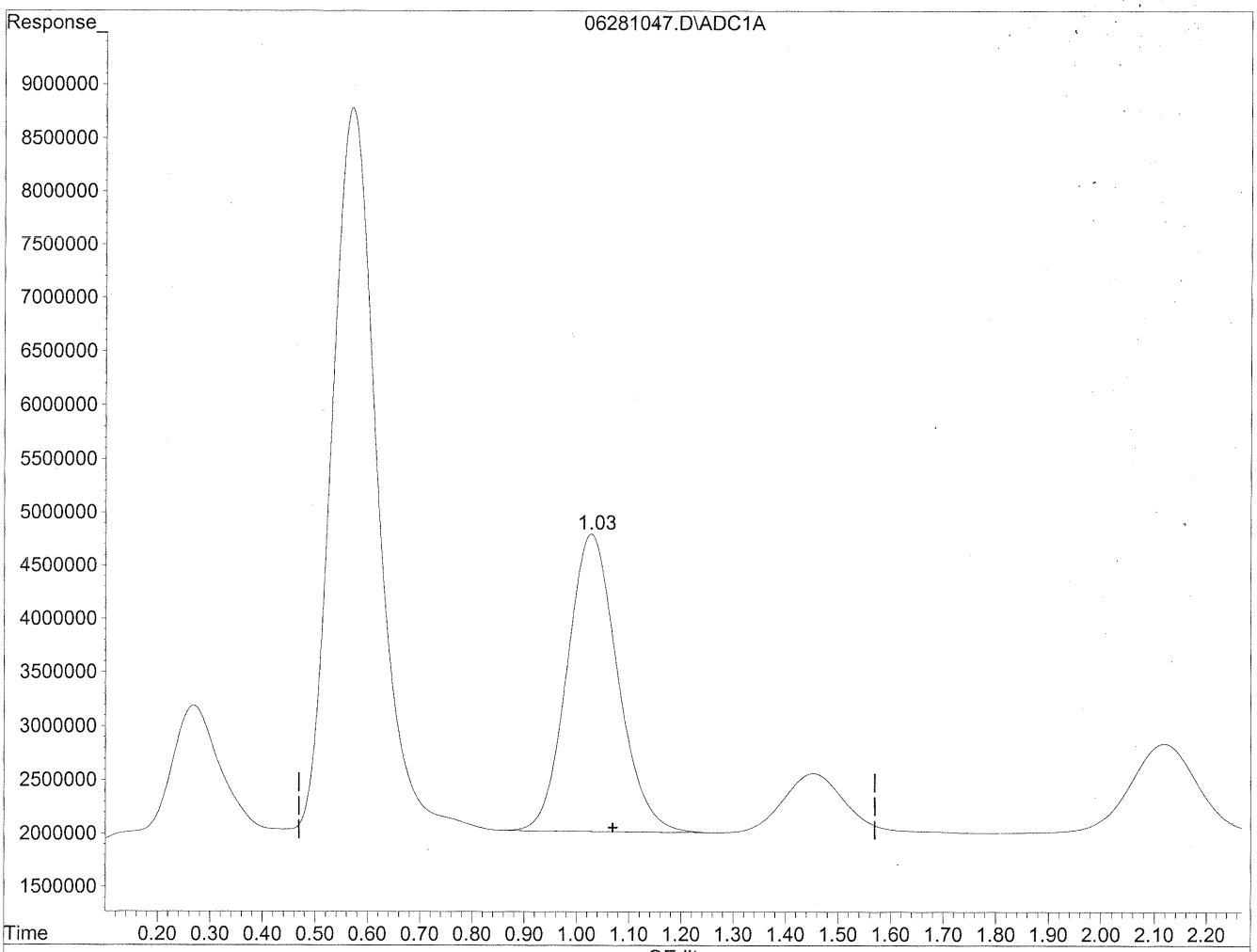
(+) = Expected Retention Time

06281047.D TO110610.M Tue Jun 29 14:56:53 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281047.D Vial: 47
Acq On : 28 Jun 2010 20:43 Operator: MD
Sample : P1002127-008 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 14:56 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.03min 1300.563ng/ml m

response 186221979

TC 6/30/10

TC 6/30/10

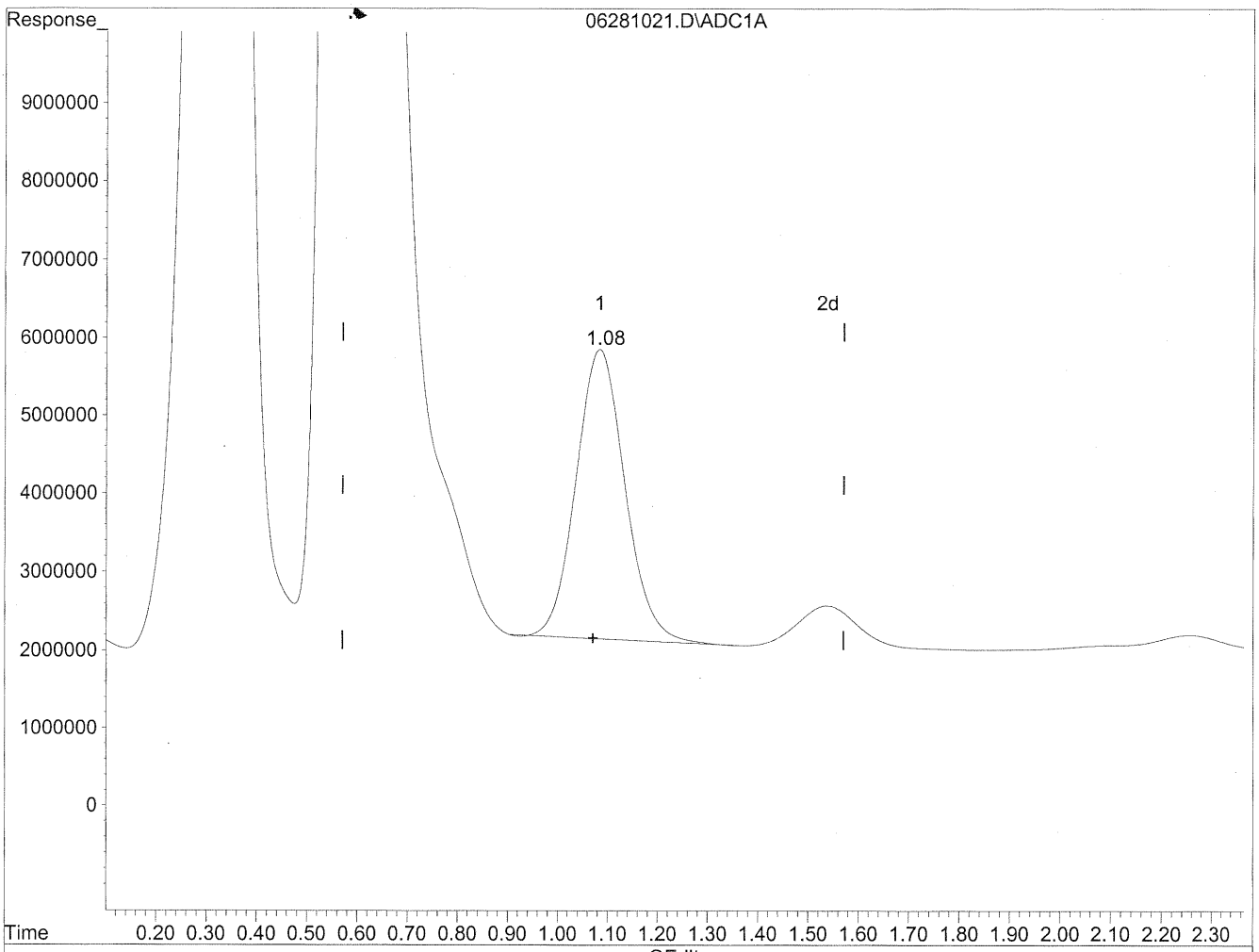
(+) = Expected Retention Time

06281047.D TO110610.M Tue Jun 29 14:57:03 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281021.D Vial: 21
Acq On : 28 Jun 2010 14:51 Operator: MD
Sample : P1002127-009 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 28 15:05 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 1753.683ng/ml

response 251102250

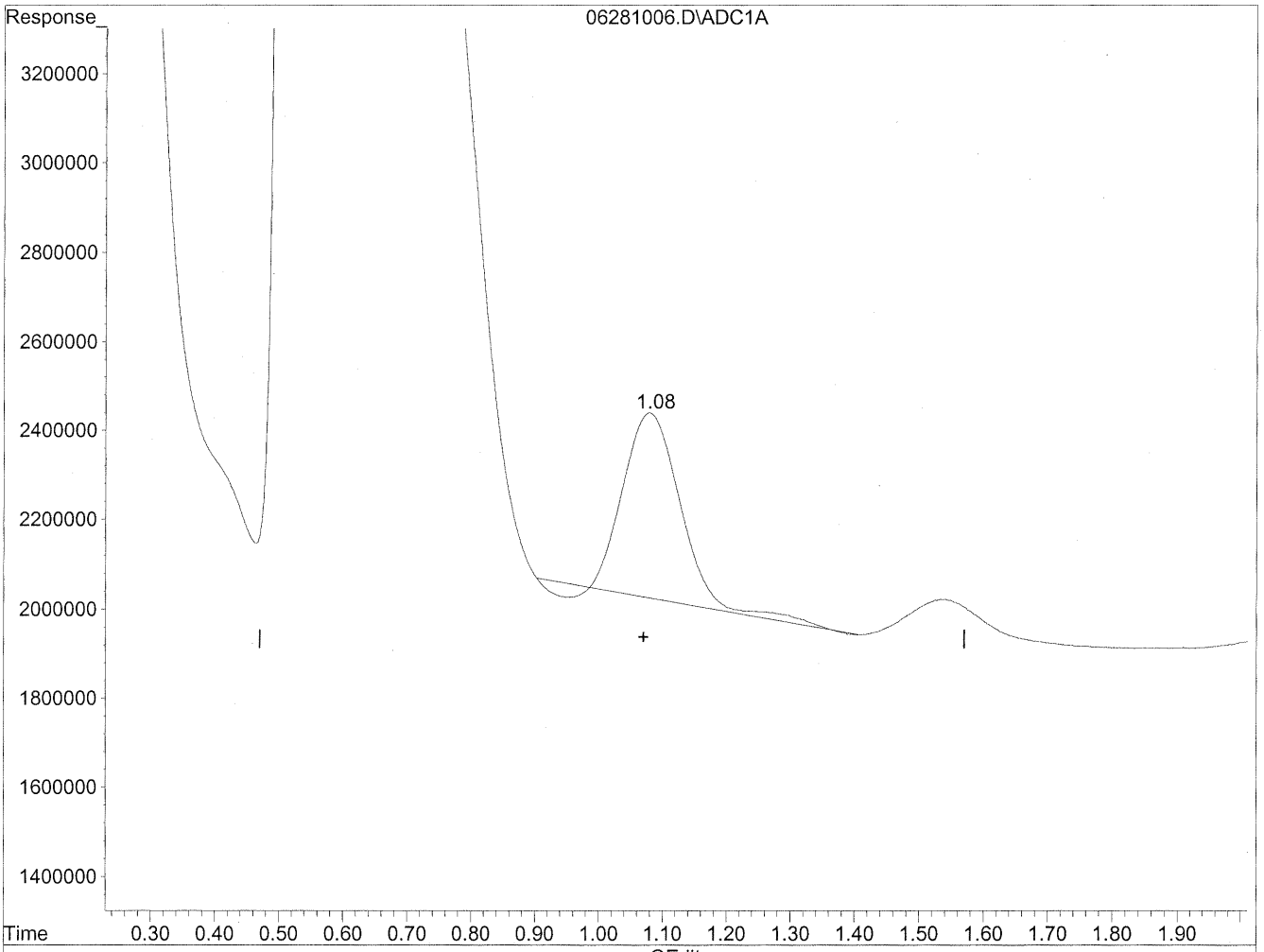
(+) = Expected Retention Time

06281021.D TO110610.M Tue Jun 29 13:24:07 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281006.D Vial: 6
Acq On : 28 Jun 2010 11:14 Operator: MD
Sample : P1002127-010 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:20 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 174.744ng/ml

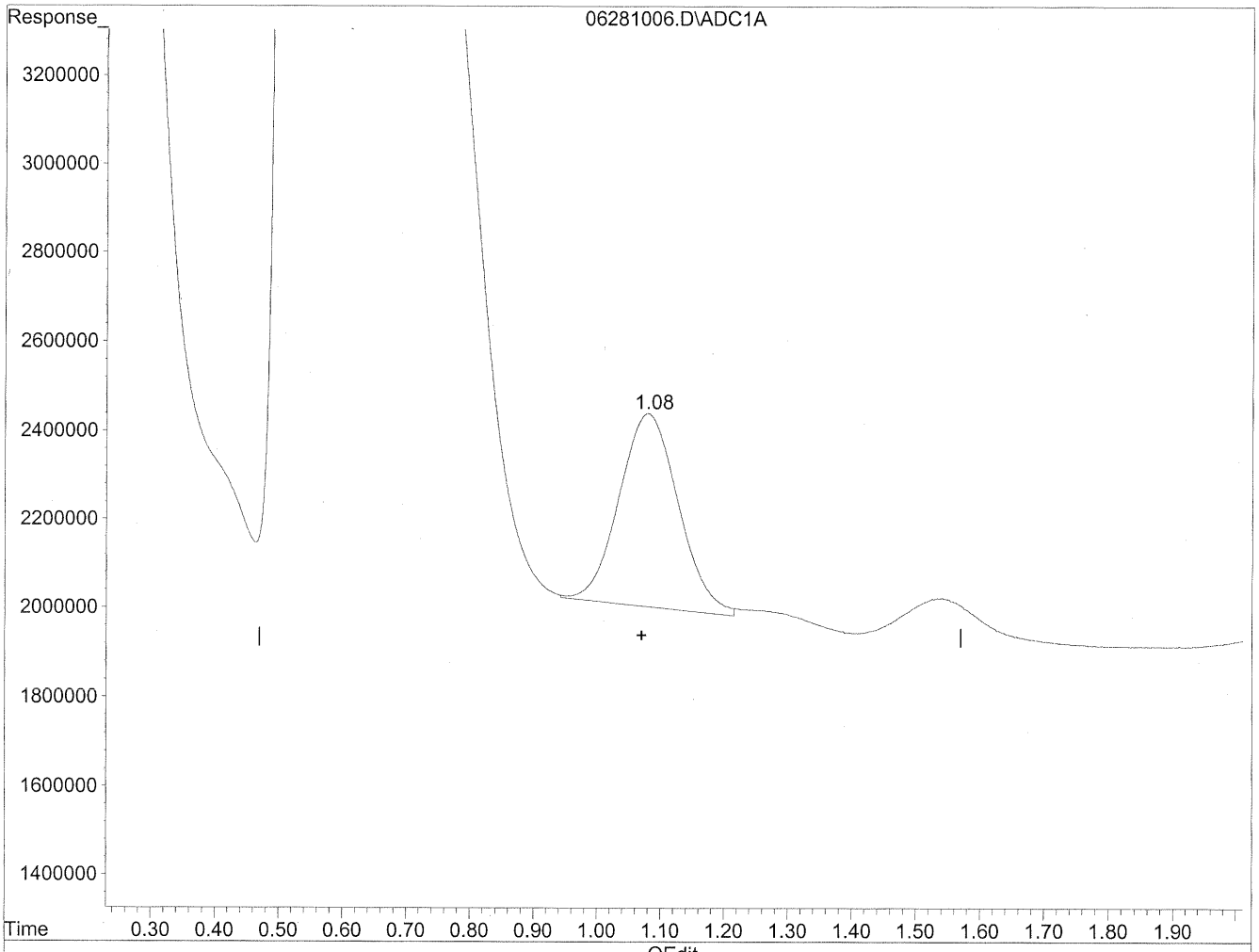
response 25020886

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281006.D Vial: 6
Acq On : 28 Jun 2010 11:14 Operator: MD
Sample : P1002127-010 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:20 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 197.629ng/ml m

response 28297592

6/29/10

6/30/10

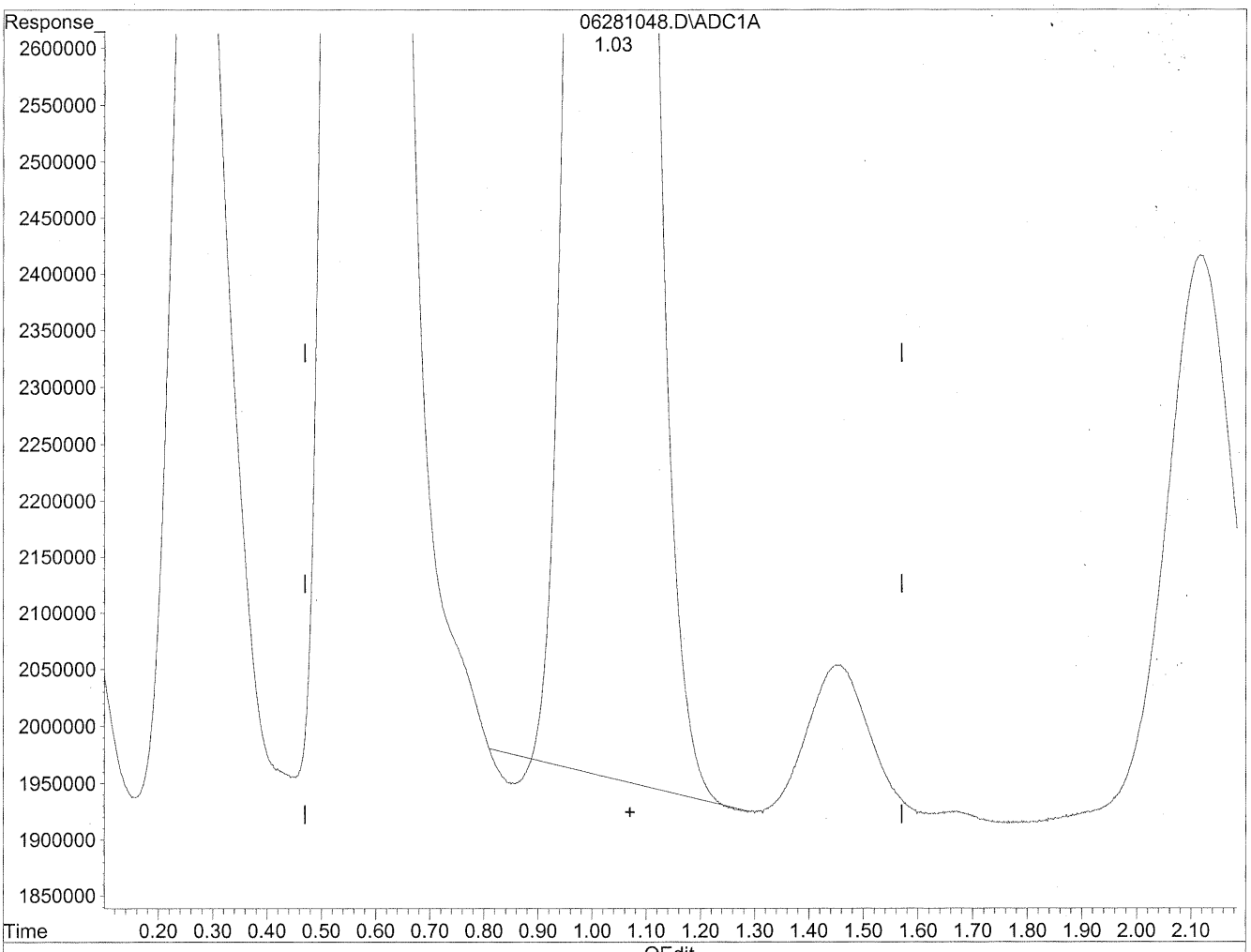
25

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281048.D Vial: 48
Acq On : 28 Jun 2010 20:57 Operator: MD
Sample : P1002127-011 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 14:57 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.03min 2195.107ng/ml

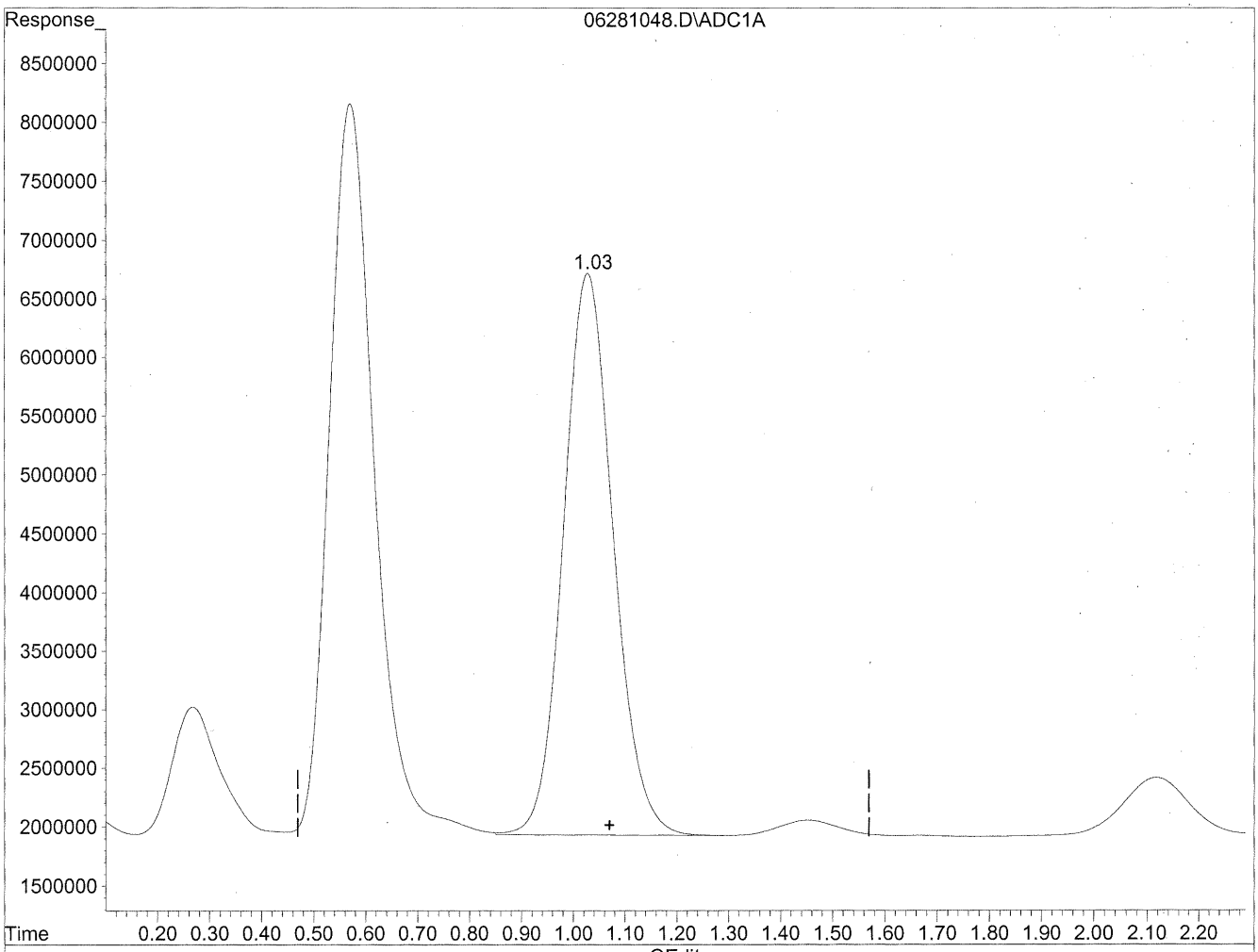
response 314307895

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281048.D Vial: 48
Acq On : 28 Jun 2010 20:57 Operator: MD
Sample : P1002127-011 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 14:57 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.03min 2237.143ng/ml m

response 320326753

6/29/10 6/30/10

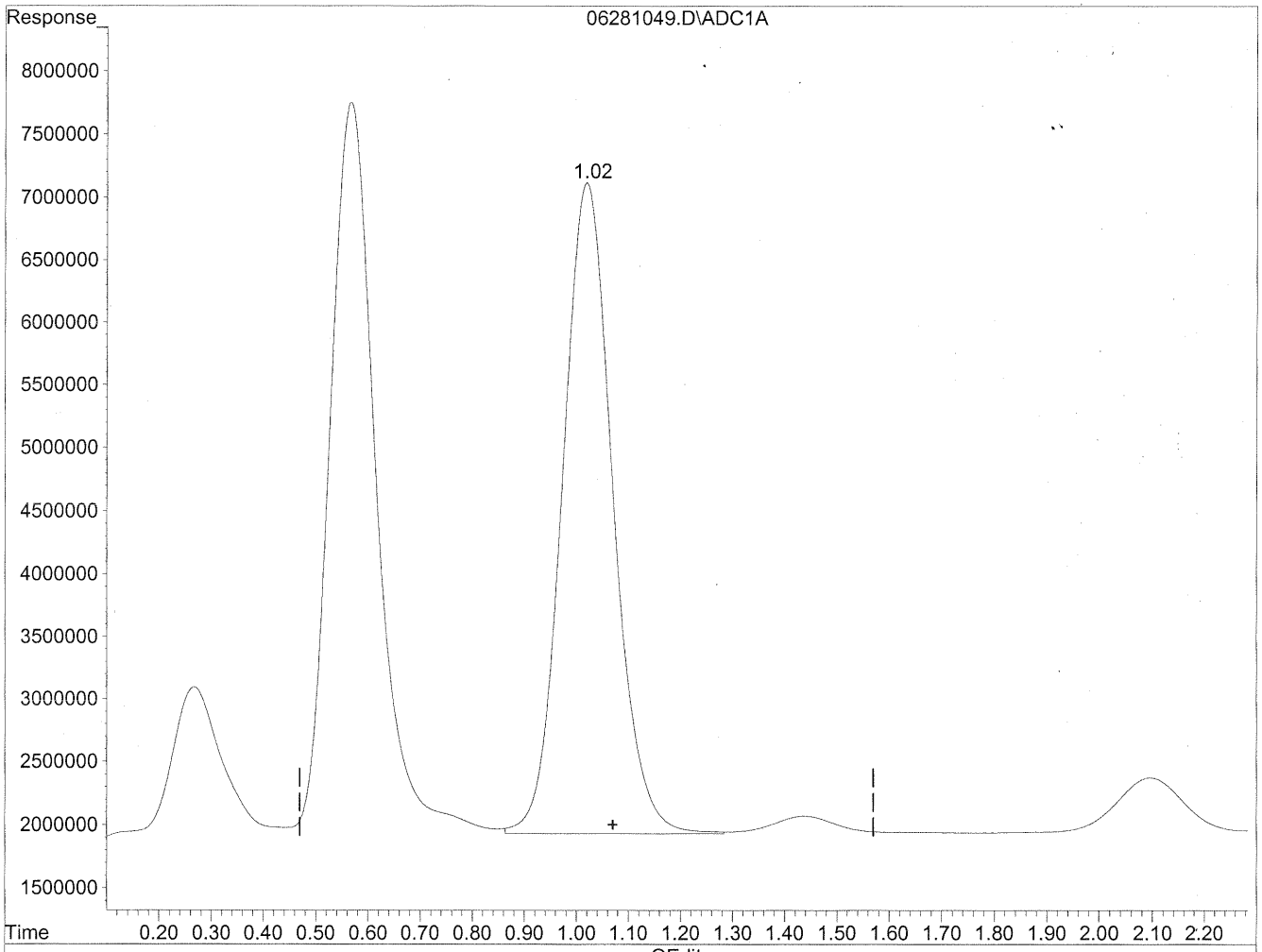
(+) = Expected Retention Time

06281048.D TO110610.M Tue Jun 29 14:57:31 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281049.D Vial: 49
Acq On : 28 Jun 2010 21:10 Operator: MD
Sample : P1002127-012 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 14:57 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.02min 2441.305ng/ml

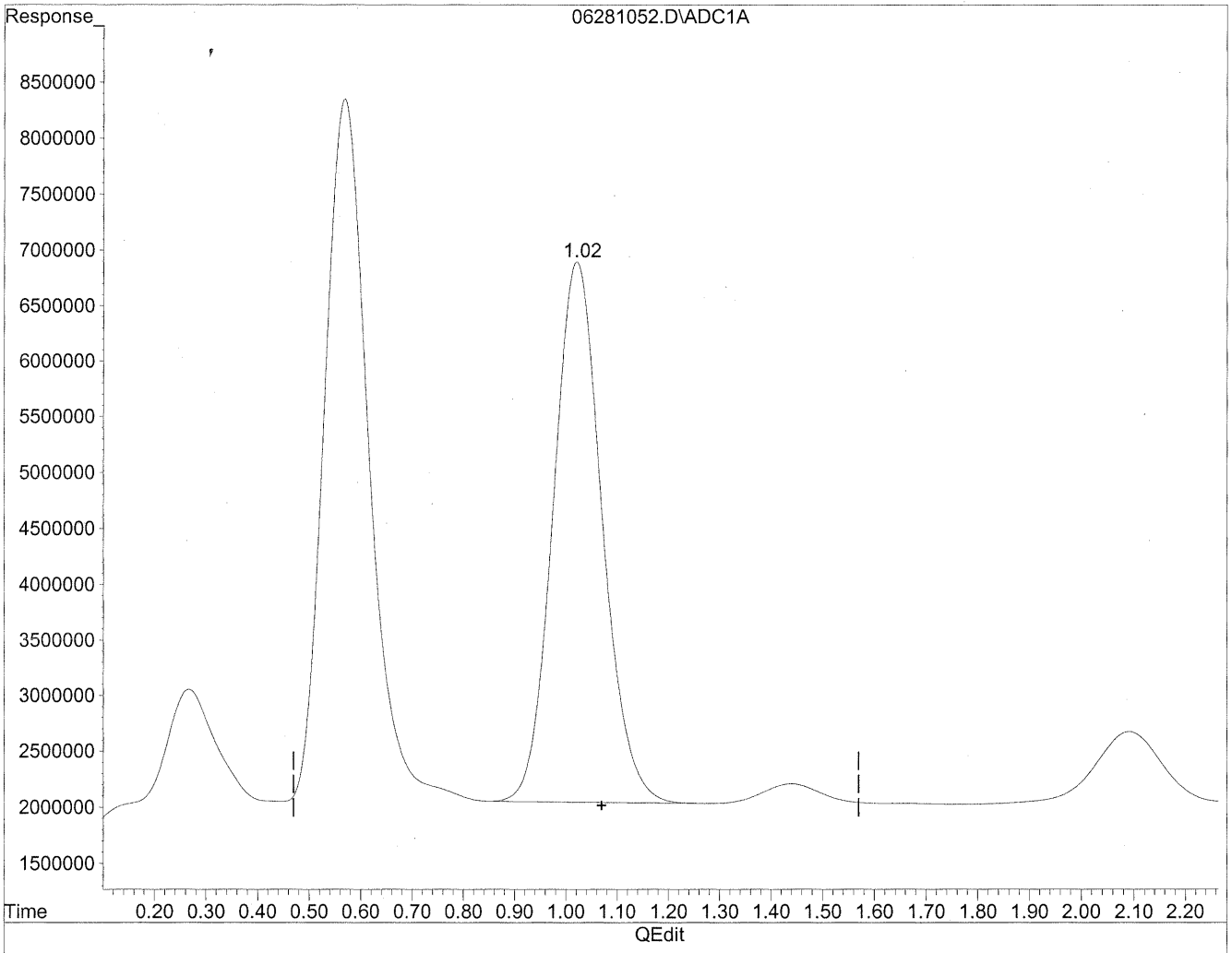
response 349559880

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281052.D Vial: 52
Acq On : 28 Jun 2010 21:51 Operator: MD
Sample : P1002127-013 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:05 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.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



(1) Formaldehyde

1.02min 2240.604ng/ml

response 320822410

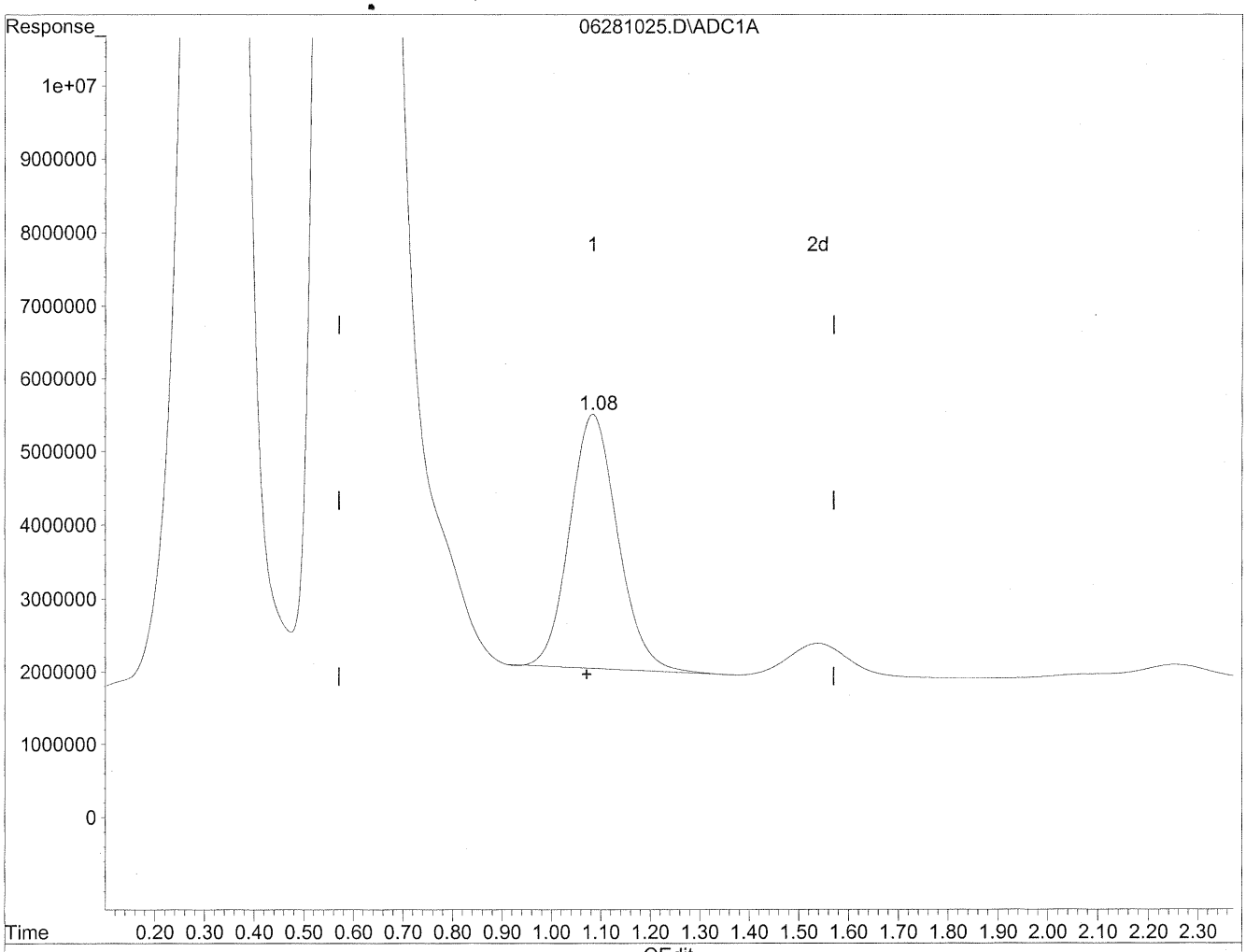
(+) = Expected Retention Time

06281052.D TO110610.M Wed Jun 30 09:25:40 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281025.D Vial: 25
Acq On : 28 Jun 2010 15:45 Operator: MD
Sample : P1002127-014 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 28 16:01 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 1638.350ng/ml

response 234588220

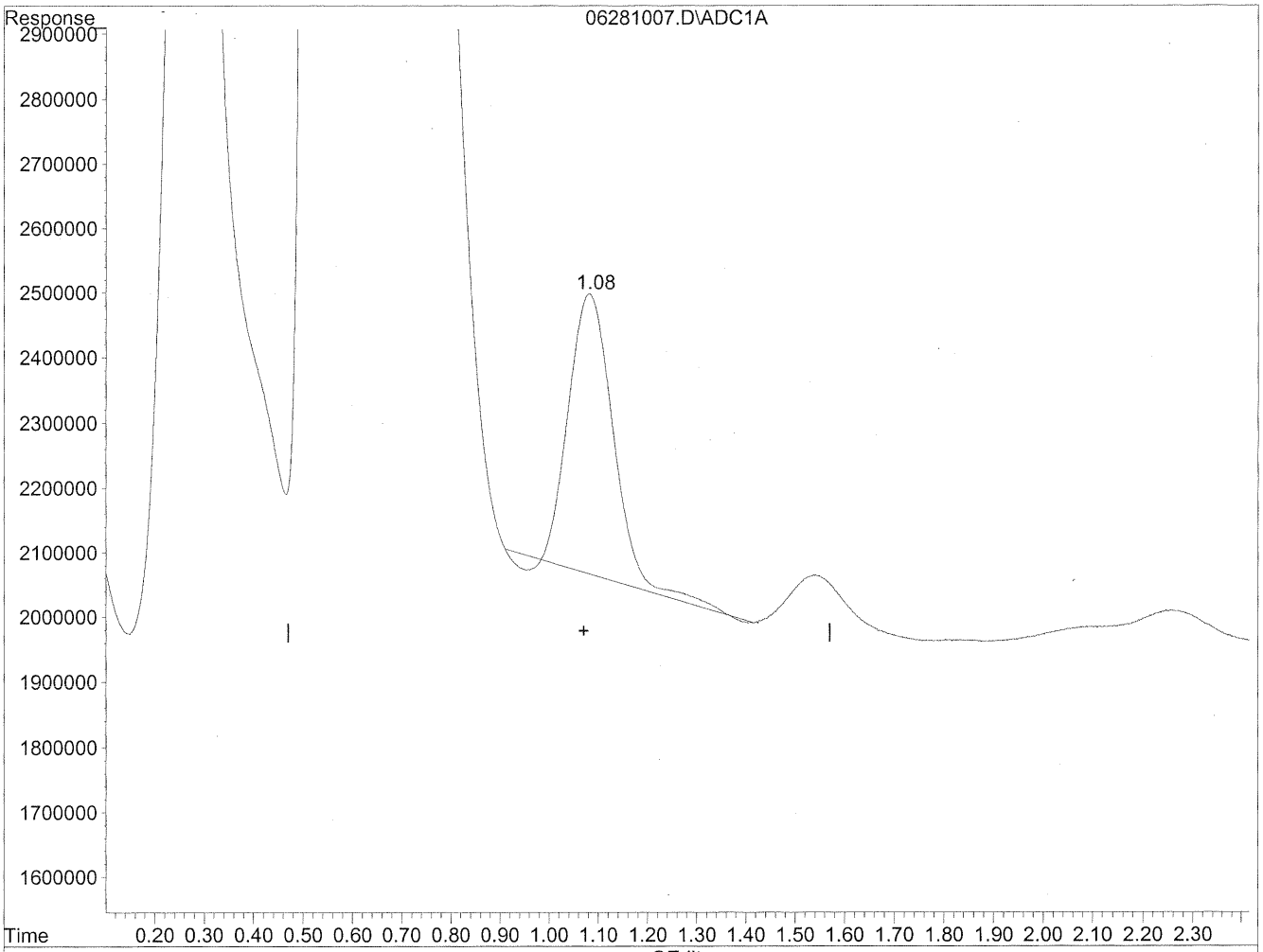
(+) = Expected Retention Time

06281025.D TO110610.M Tue Jun 29 13:24:47 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281007.D Vial: 7
Acq On : 28 Jun 2010 11:28 Operator: MD
Sample : P1002127-015 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:20 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.08min 186.558ng/ml

response 26712369

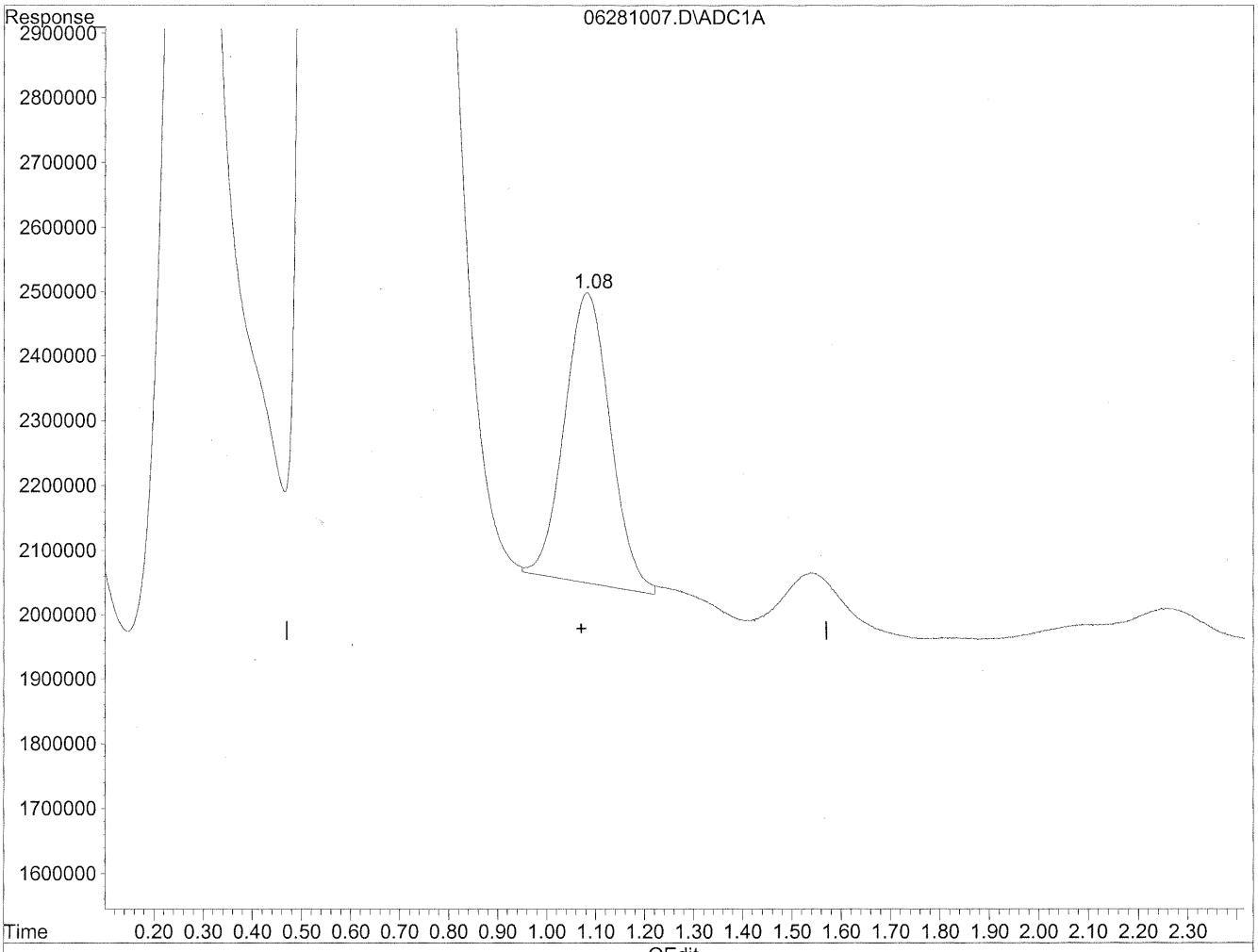
(+) = Expected Retention Time

06281007.D TO110610.M Tue Jun 29 13:21:15 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281007.D Vial: 7
Acq On : 28 Jun 2010 11:28 Operator: MD
Sample : P1002127-015 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:20 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 203.591ng/ml m

response 29151333

TC 6/30/10

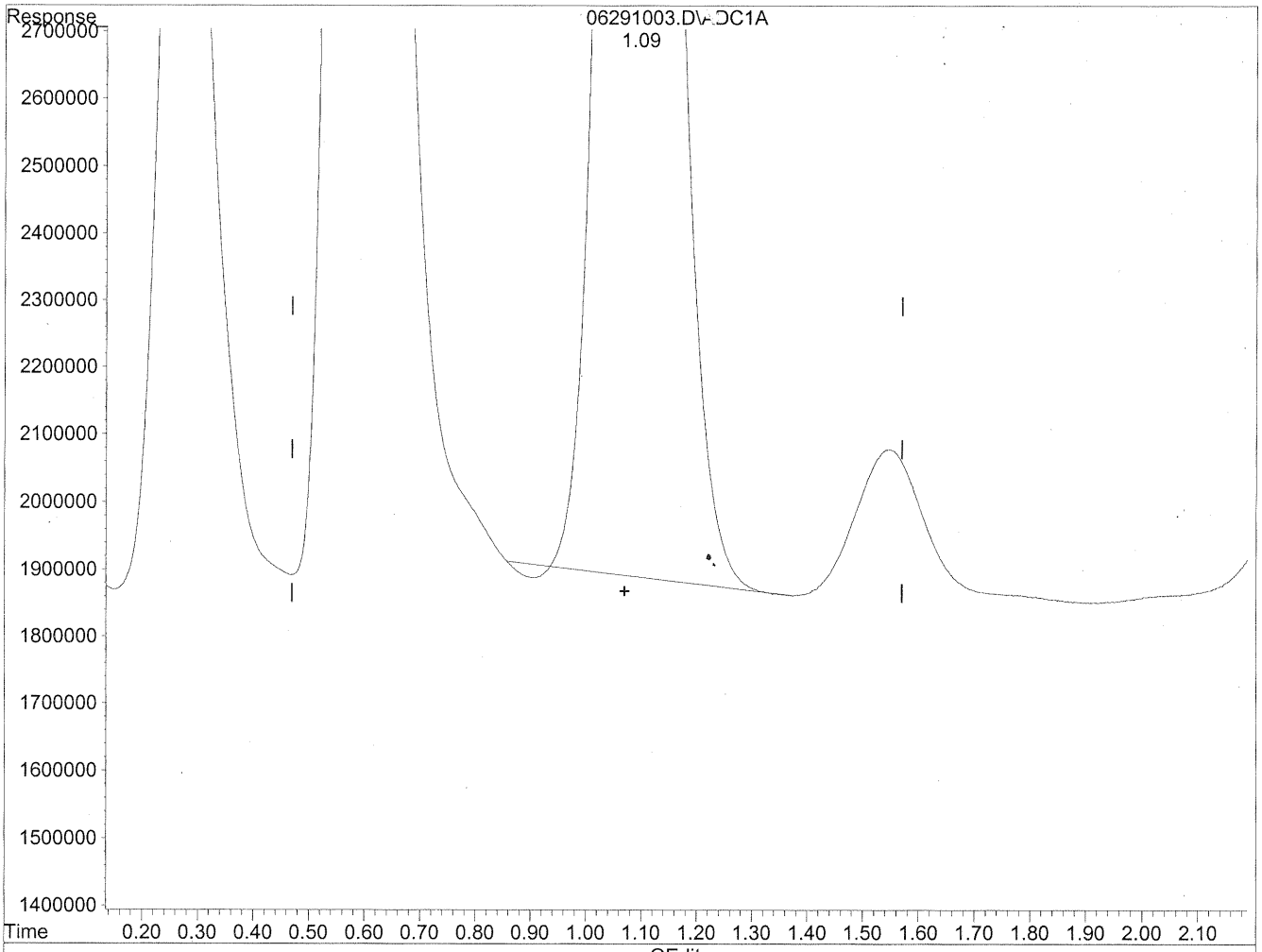
32

(+) = Expected Retention Time
06281007.D TO110610.M Tue Jun 29 13:21:28 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291003.D Vial: 3
Acq On : 29 Jun 2010 10:19 Operator: MD
Sample : P1002127-016 2.0ml 10x dil Inst : LC 01
Misc : Radiello passive sampler Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:08 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.09min 2139.605ng/ml

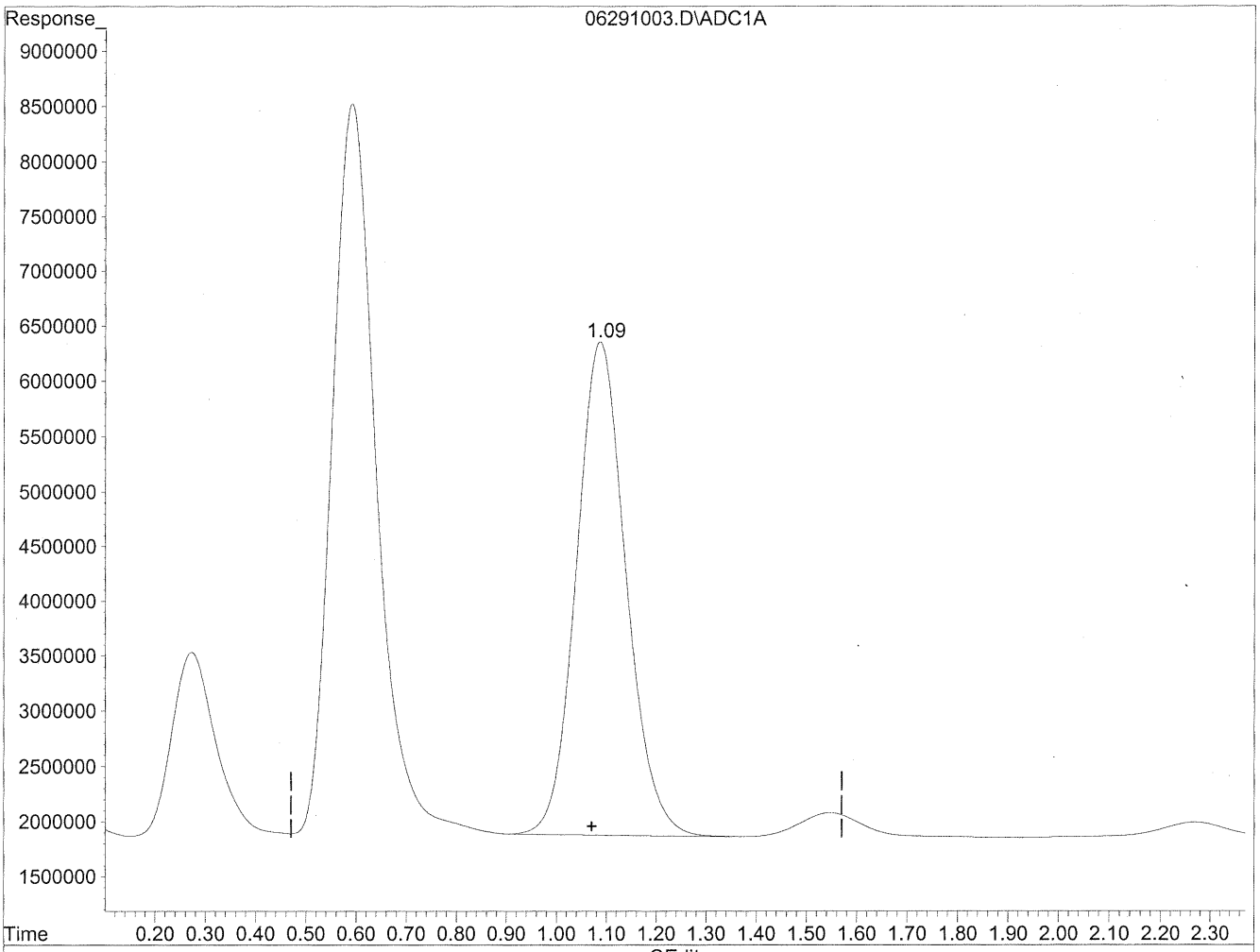
response 306360810

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291003.D Vial: 3
Acq On : 29 Jun 2010 10:19 Operator: MD
Sample : P1002127-016 2.0ml 10x dil Inst : LC 01
Misc : Radiello passive sampler Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:08 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.09min 2163.027ng/ml m

response 309714407

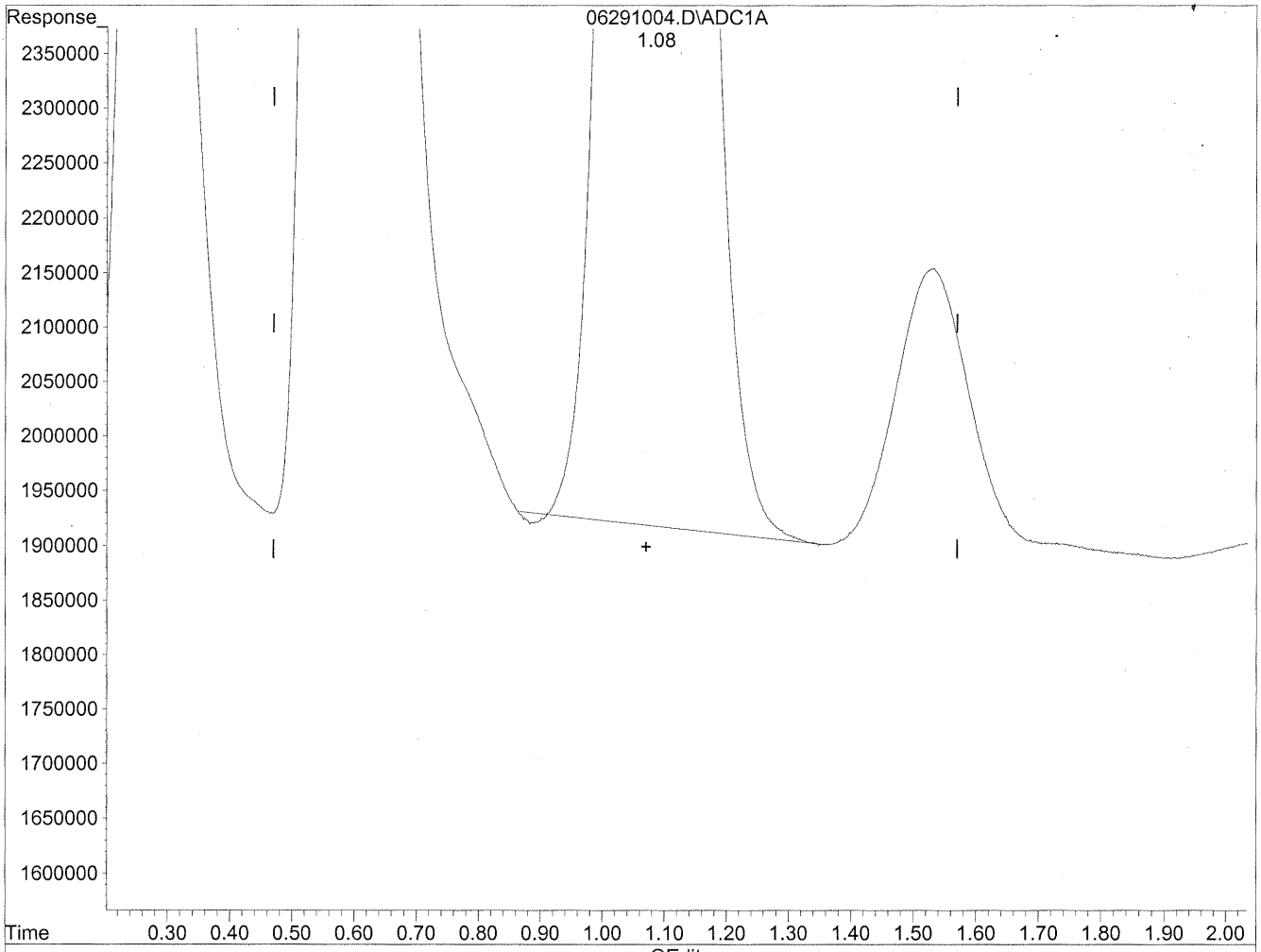
6/30/10 JC 6/30/10

(+) = Expected Retention Time
06291003.D TO110610.M Tue Jun 29 15:08:24 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291004.D Vial: 4
Acq On : 29 Jun 2010 10:32 Operator: MD
Sample : P1002127-017 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:08 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.08min 2387.219ng/ml

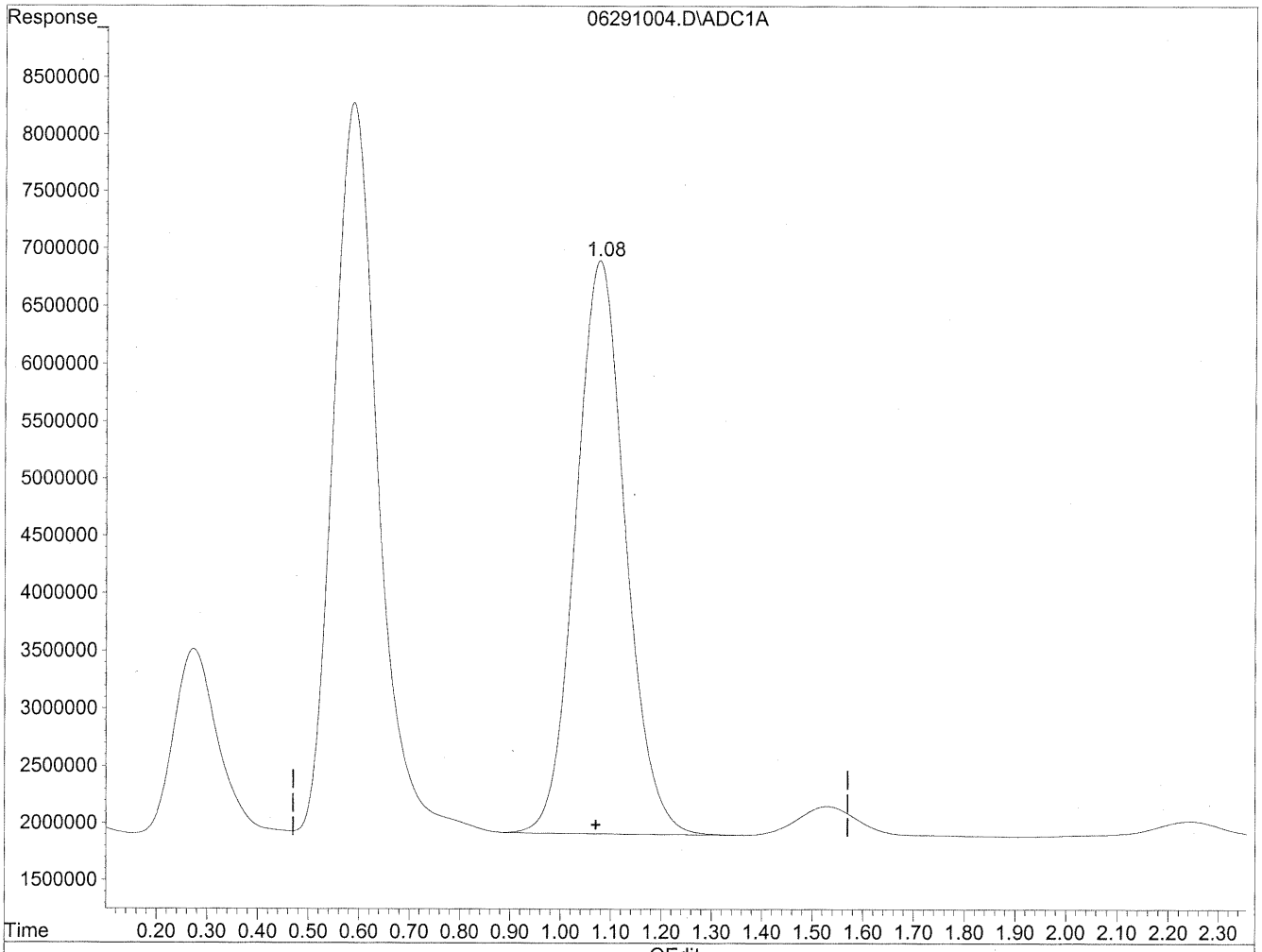
response 341815584

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291004.D Vial: 4
Acq On : 29 Jun 2010 10:32 Operator: MD
Sample : P1002127-017 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:08 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 2399.079ng/ml m

response 343513667

MC 6/30/10 TC 6/30/10

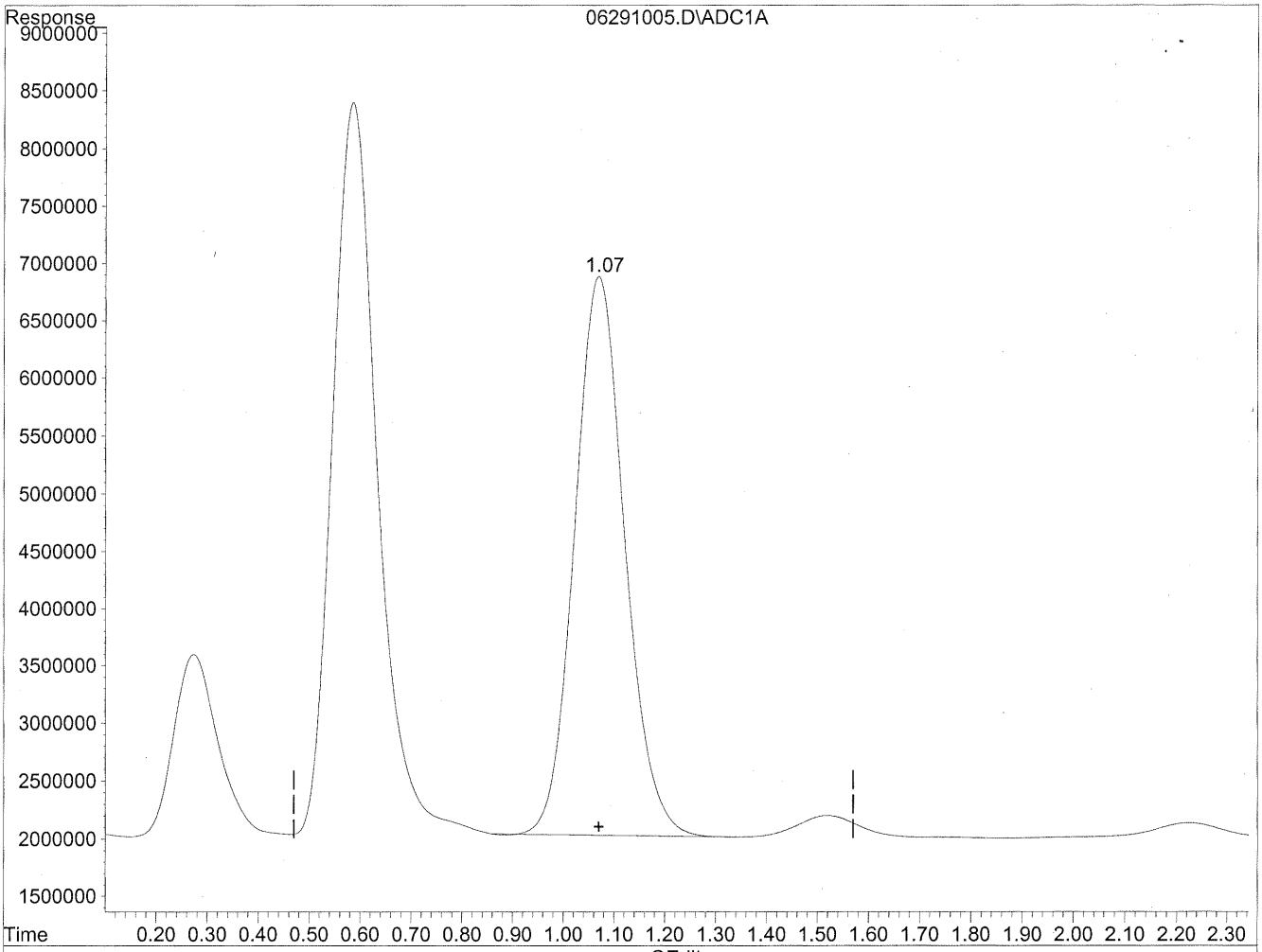
(+) = Expected Retention Time

06291004.D TO110610.M Tue Jun 29 15:08:40 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291005.D Vial: 5
Acq On : 29 Jun 2010 10:46 Operator: MD
Sample : P1002127-018 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:08 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.07min 2326.087ng/ml

response 333062319

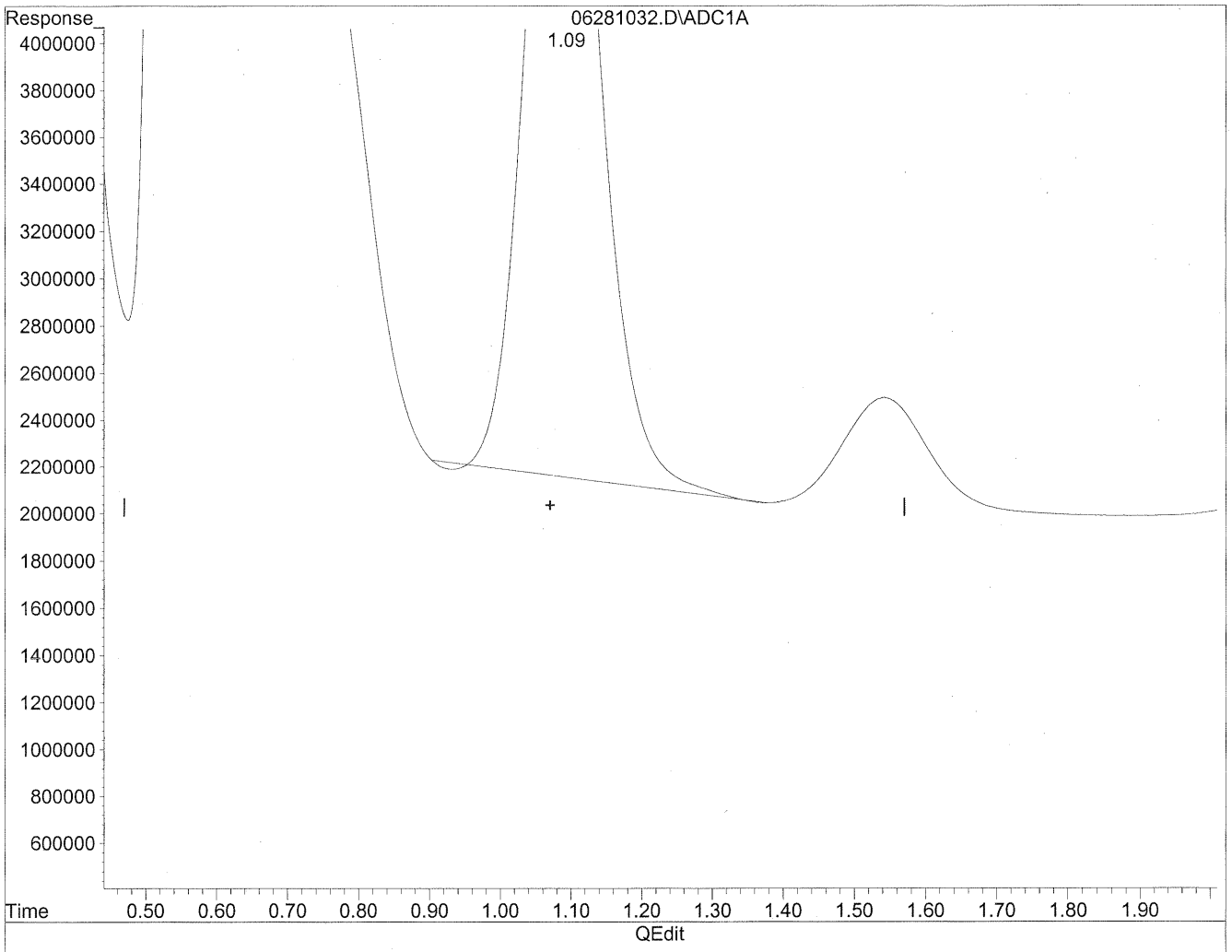
(+) = Expected Retention Time

06291005.D TO110610.M Tue Jun 29 15:08:51 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281032.D Vial: 32
Acq On : 28 Jun 2010 17:20 Operator: MD
Sample : P1002127-019 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:25 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.09min 1788.958ng/ml

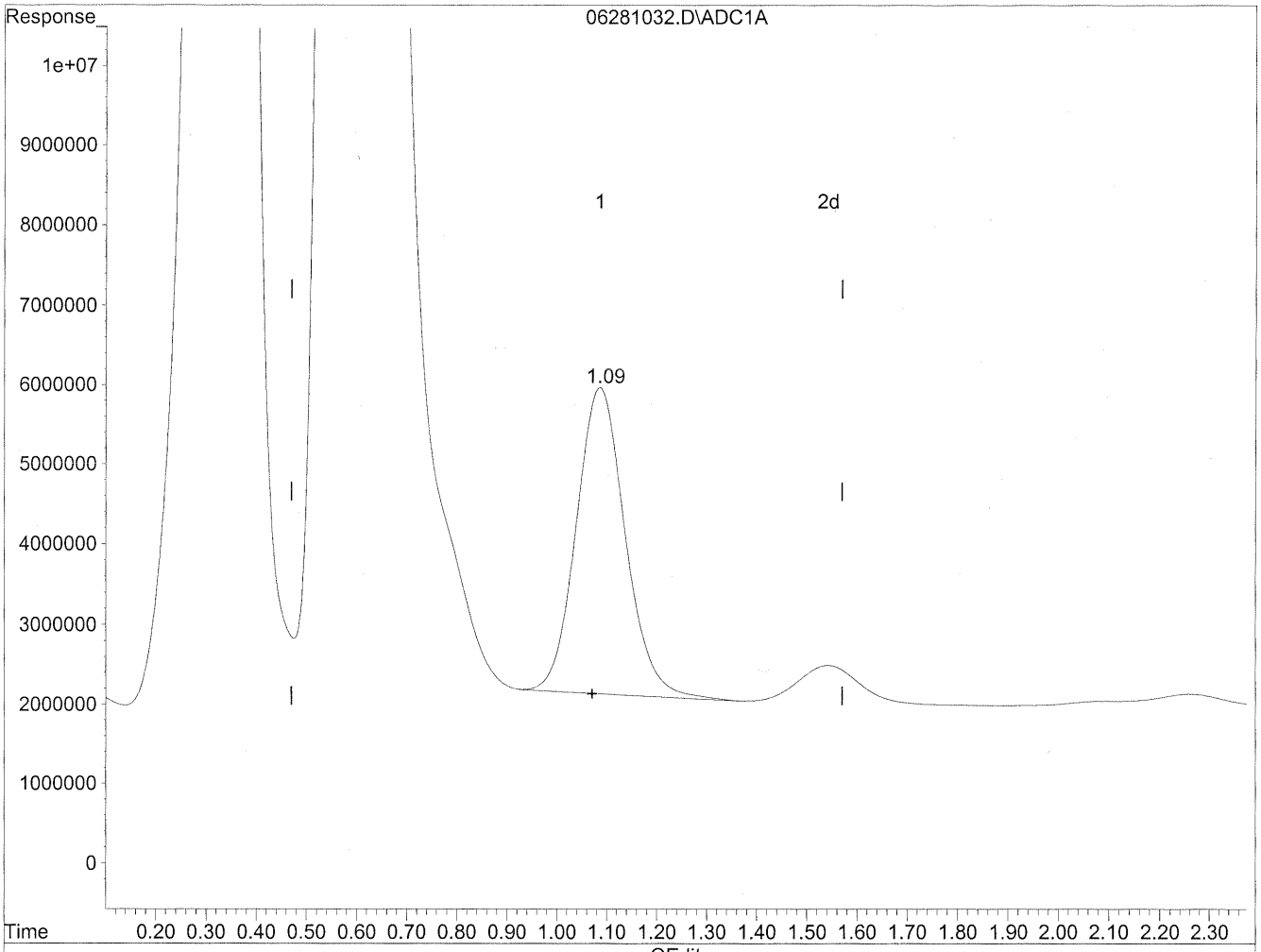
response 256153081

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281032.D Vial: 32
Acq On : 28 Jun 2010 17:20 Operator: MD
Sample : P1002127-019 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:25 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.09min 1828.361ng/ml m

response 261795021

6/30/10

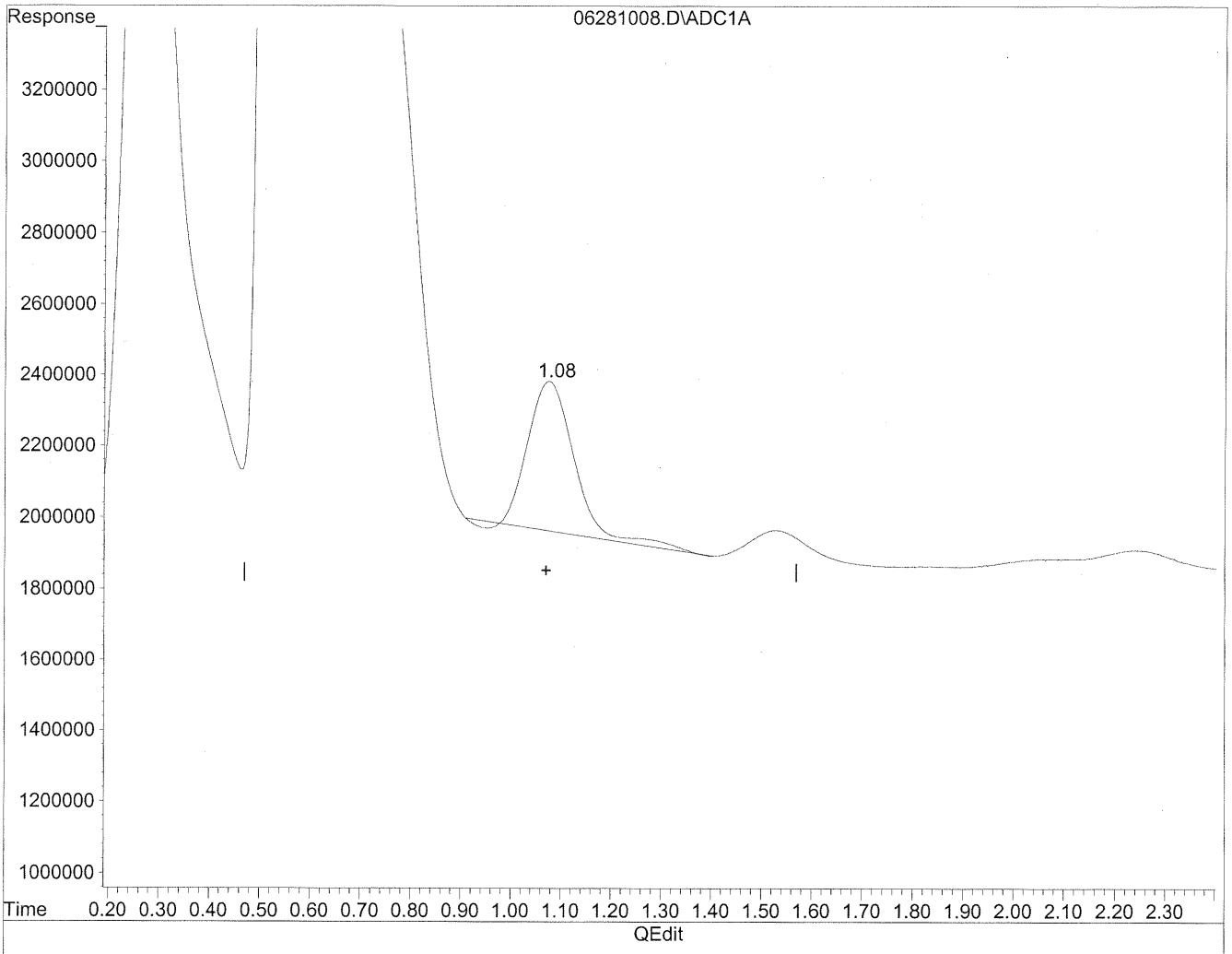
6/30/10

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281008.D Vial: 8
Acq On : 28 Jun 2010 11:42 Operator: MD
Sample : P1002127-020 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:21 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.07min 186.241ng/ml

response 26667076

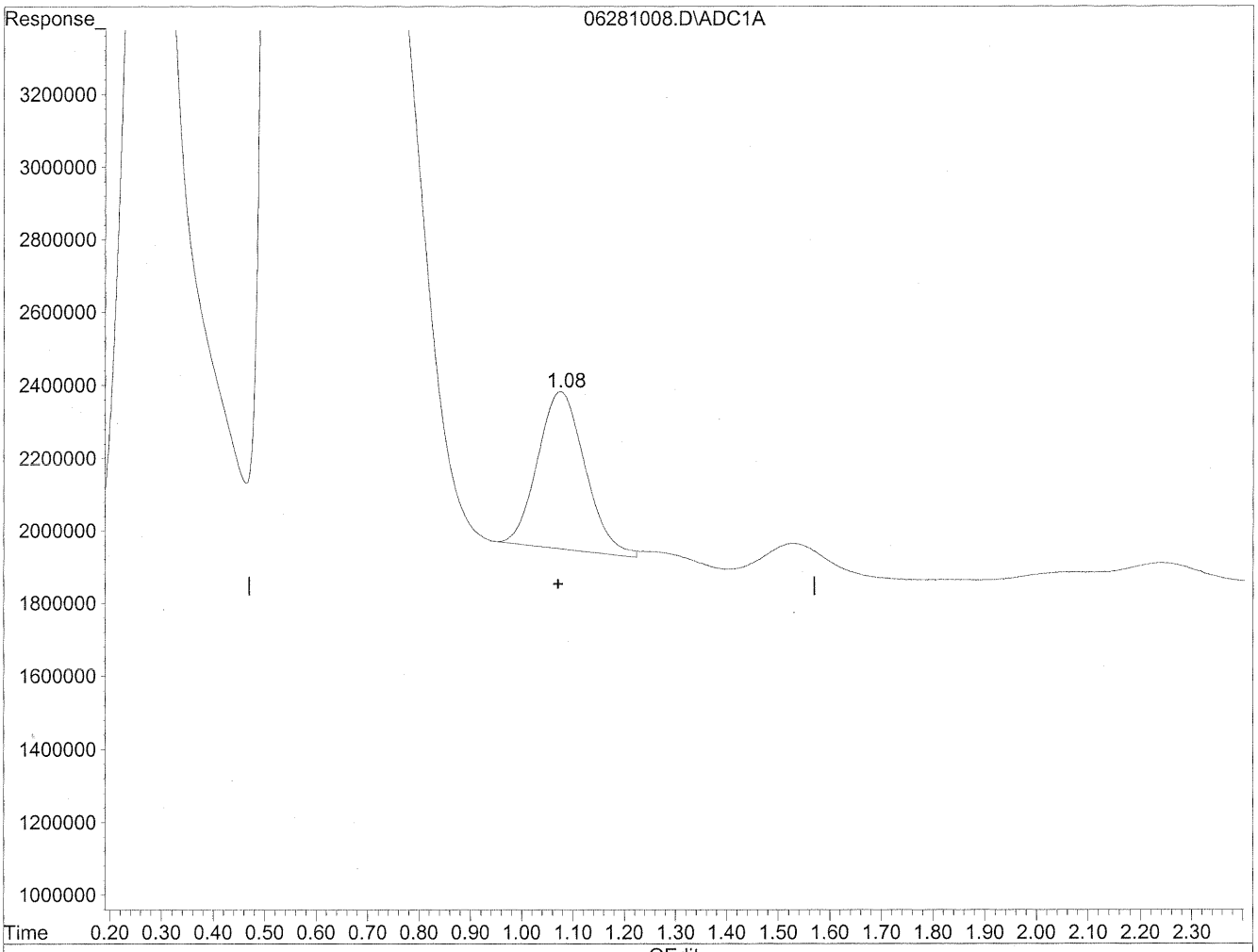
(+) = Expected Retention Time

06281008.D TO110610.M Tue Jun 29 13:21:39 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281008.D Vial: 8
Acq On : 28 Jun 2010 11:42 Operator: MD
Sample : P1002127-020 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:21 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.08min 194.682ng/ml m

response 27875742

4/30/10

6/30/10

41

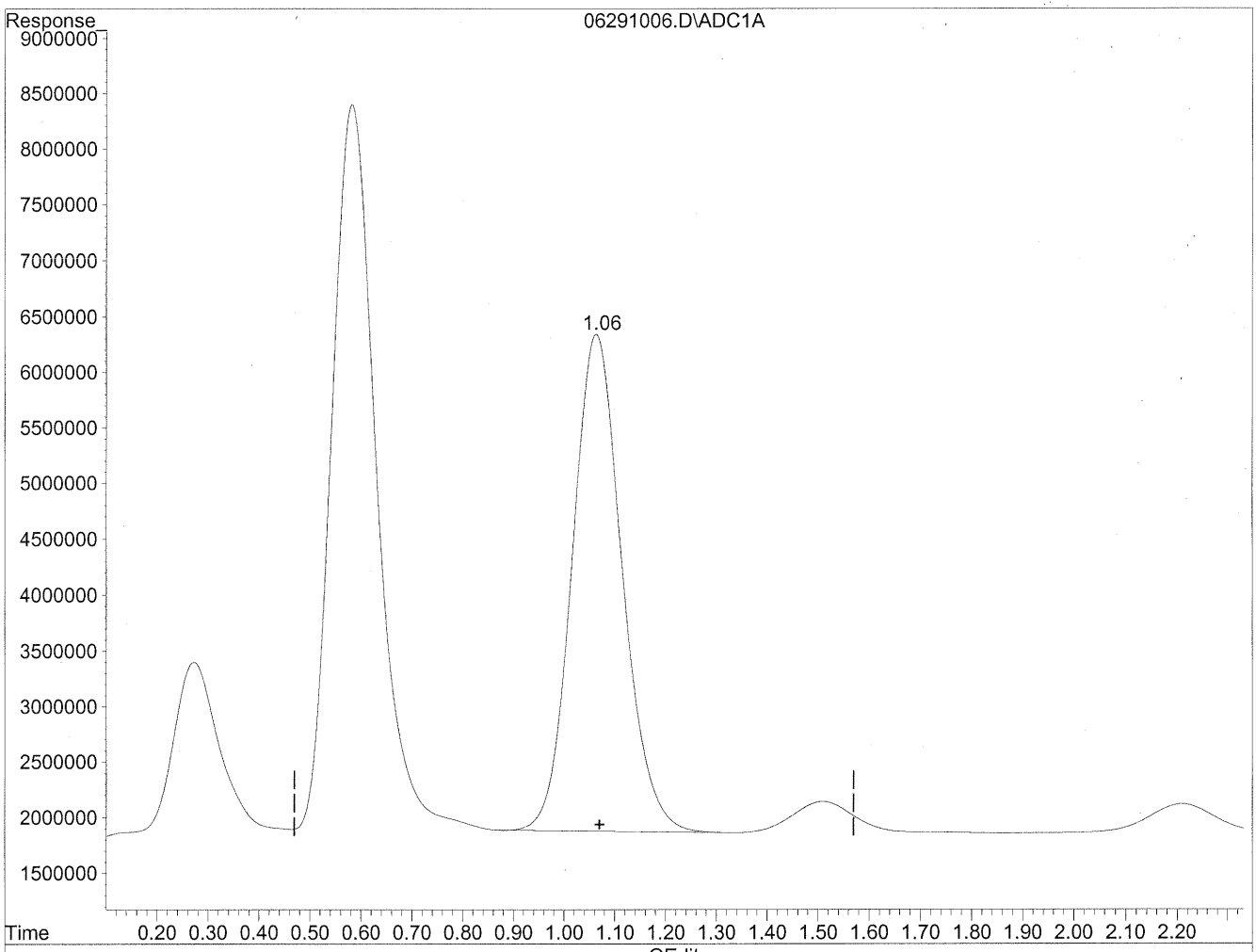
(+) = Expected Retention Time

06281008.D TO110610.M Tue Jun 29 13:21:46 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291006.D Vial: 6
Acq On : 29 Jun 2010 10:59 Operator: MD
Sample : P1002127-021 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:08 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.06min 2135.412ng/ml

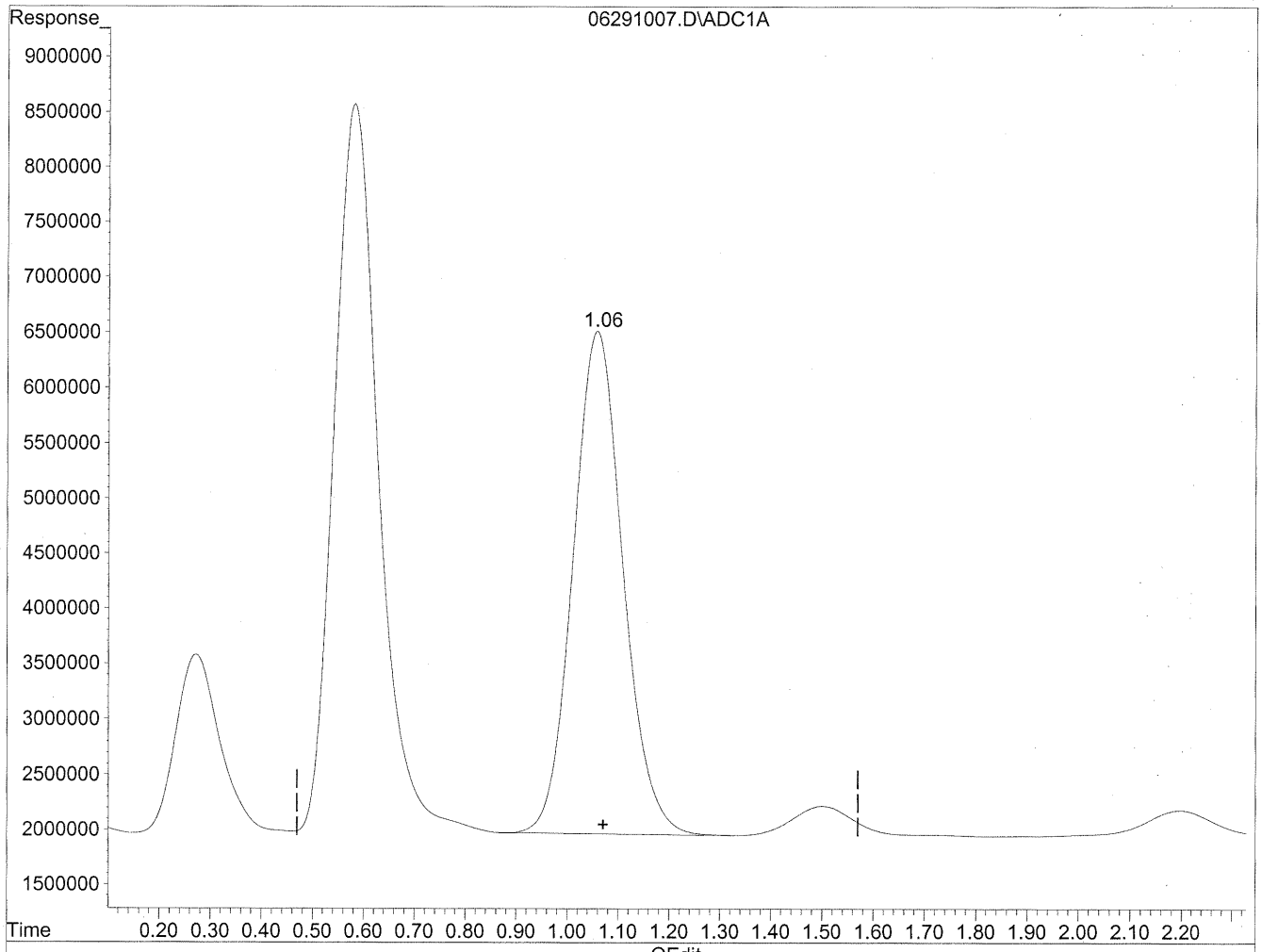
response 305760455

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291007.D Vial: 7
Acq On : 29 Jun 2010 11:13 Operator: MD
Sample : P1002127-022 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:09 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.06min 2163.404ng/ml

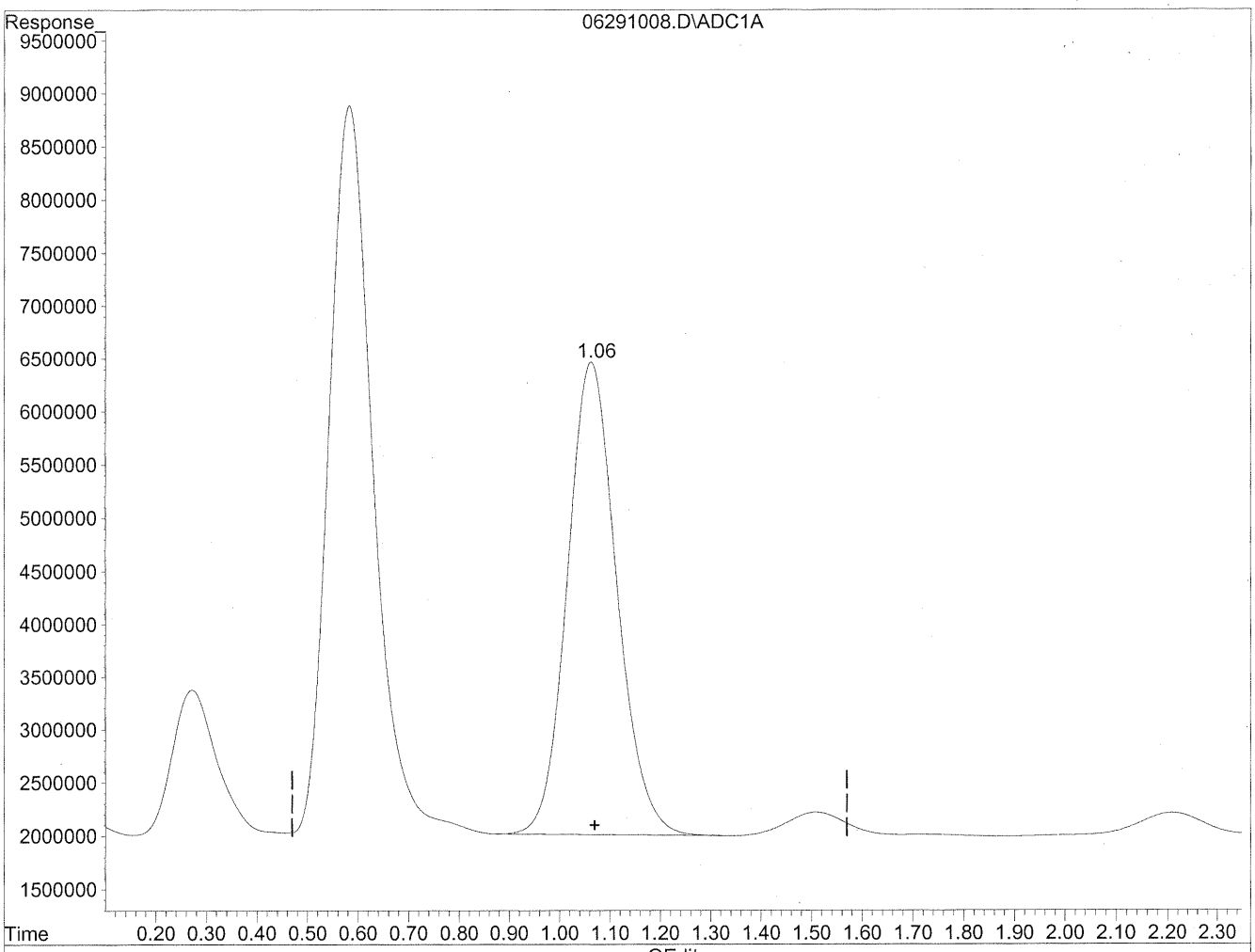
response 309768503

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291008.D Vial: 8
Acq On : 29 Jun 2010 11:27 Operator: MD
Sample : P1002127-023 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:09 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.06min 2122.313ng/ml

response 303884851

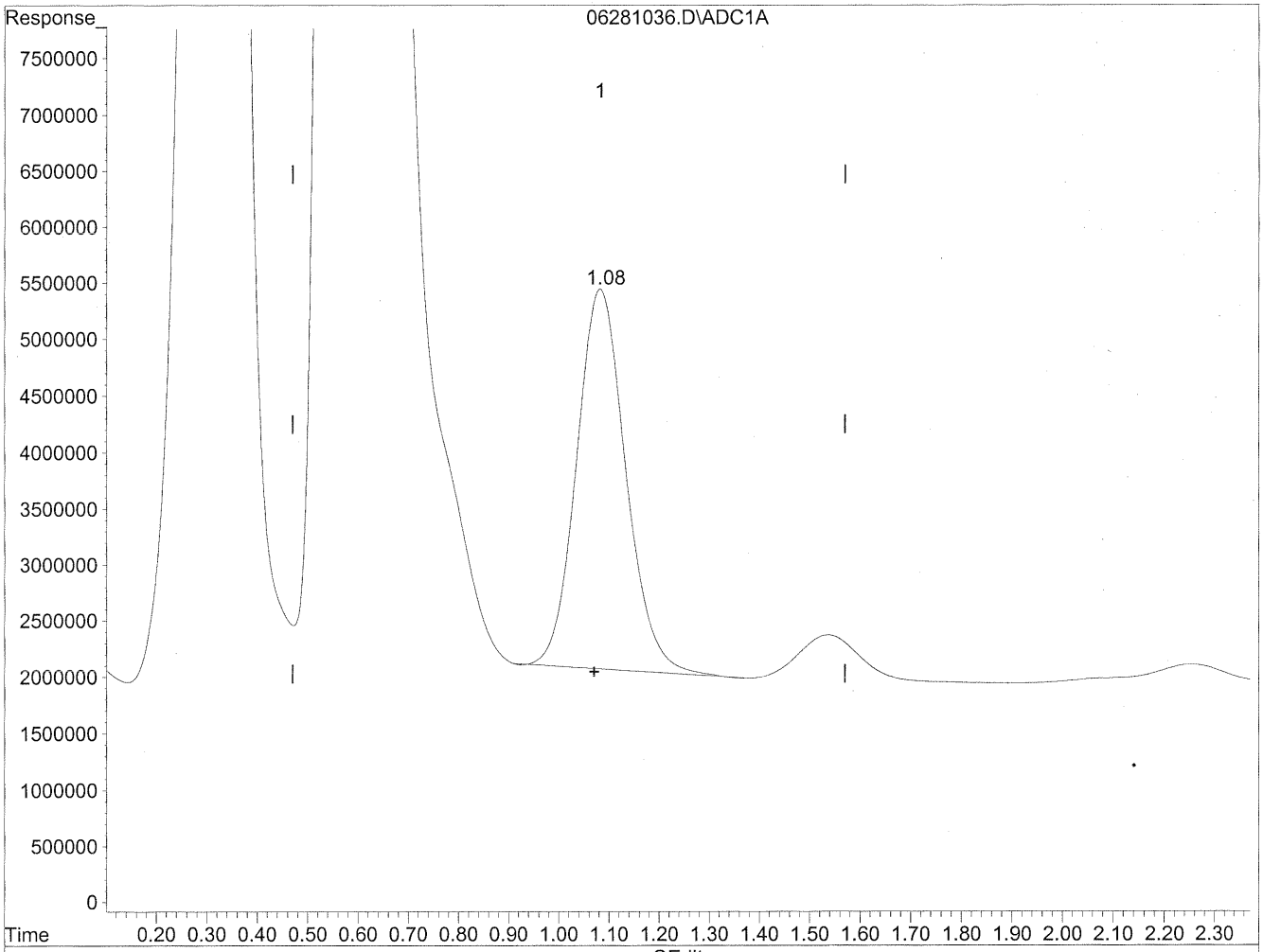
(+) = Expected Retention Time

06291008.D TO110610.M Tue Jun 29 15:09:24 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281036.D Vial: 36
Acq On : 28 Jun 2010 18:14 Operator: MD
Sample : P1002127-024 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 1606.758ng/ml

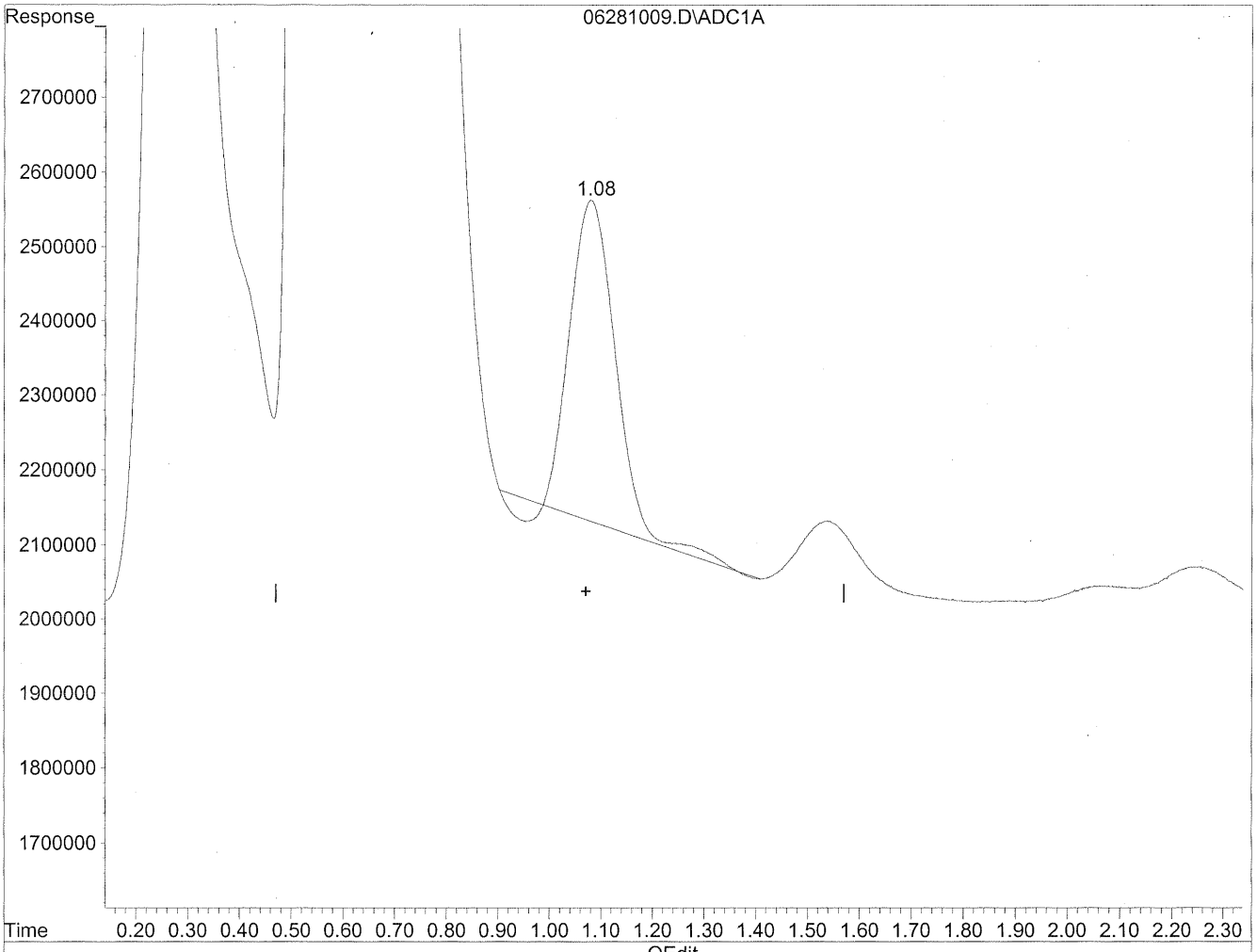
response 230064717

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281009.D Vial: 9
Acq On : 28 Jun 2010 11:55 Operator: MD
Sample : P1002127-025 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:21 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 180.360ng/ml

response 25824933

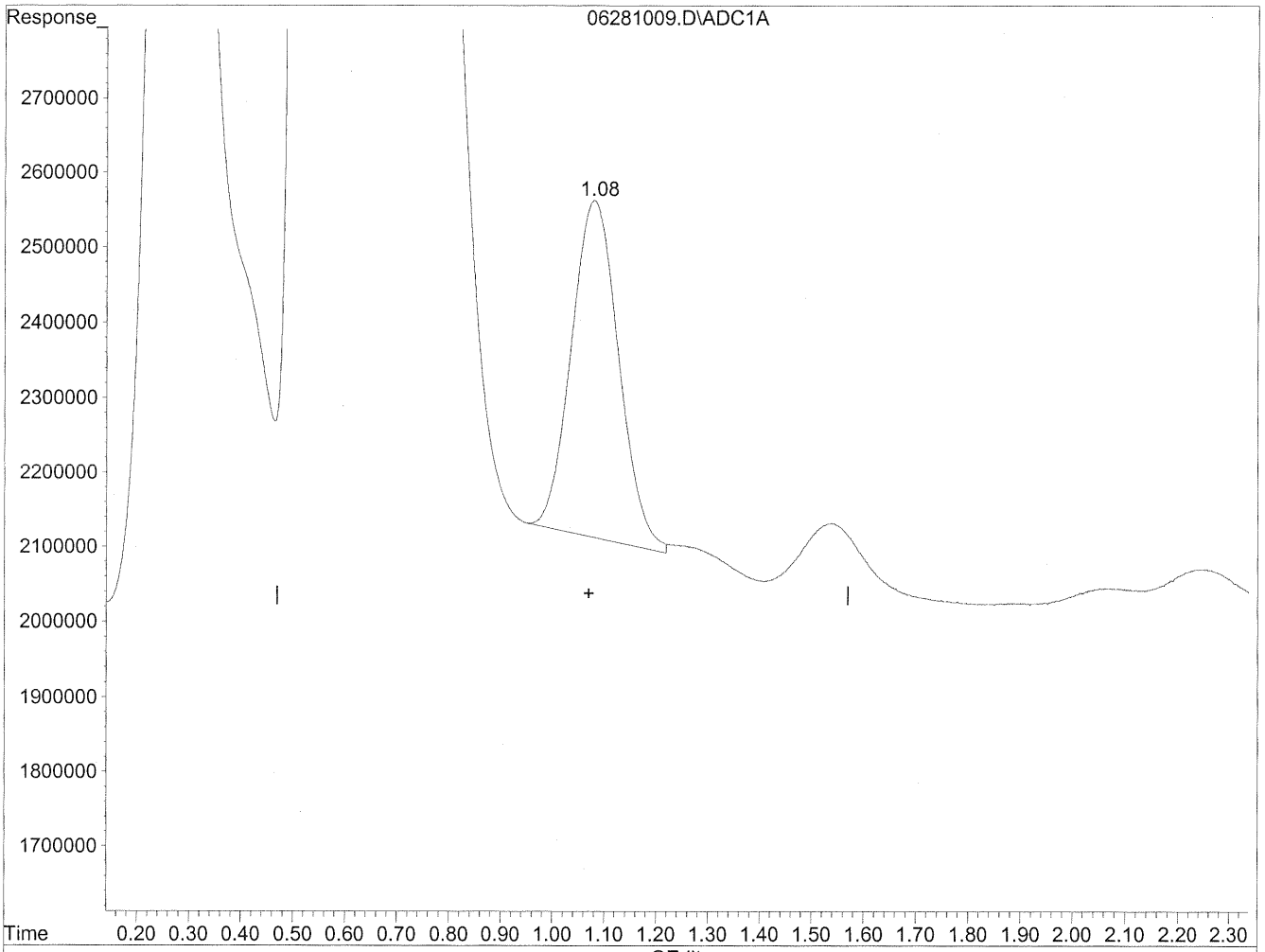
(+) = Expected Retention Time

06281009.D TO110610.M Tue Jun 29 13:21:58 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281009.D Vial: 9
Acq On : 28 Jun 2010 11:55 Operator: MD
Sample : P1002127-025 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:21 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.08min 200.166ng/ml m

response 28660897

MC 6/30/10

ME 6/30/10

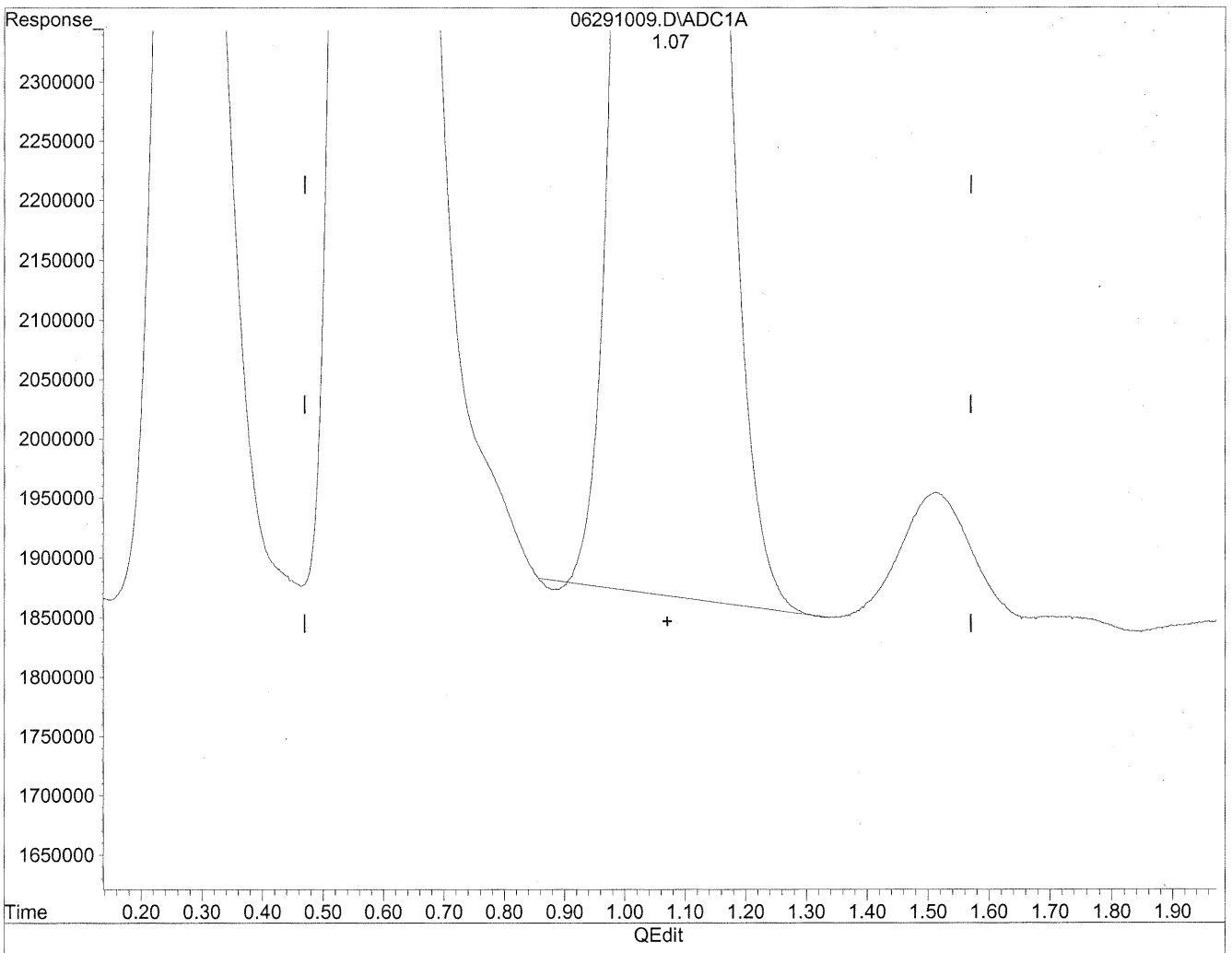
(+) = Expected Retention Time

06281009.D TO110610.M Tue Jun 29 13:22:04 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291009.D Vial: 9
Acq On : 29 Jun 2010 11:40 Operator: MD
Sample : P1002127-026 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:09 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.07min 2267.650ng/ml

response 324695030

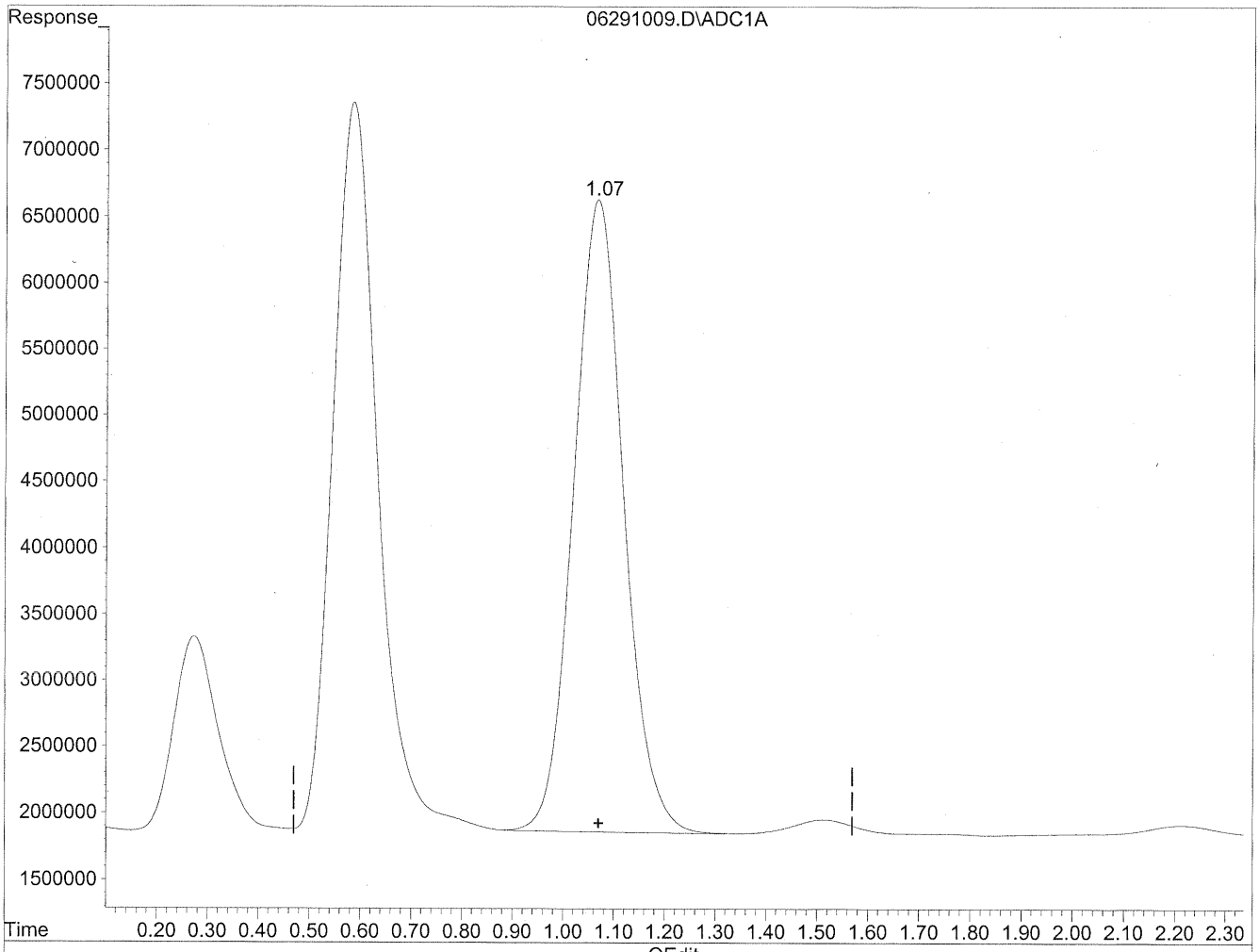
(+) = Expected Retention Time

06291009.D TO110610.M Tue Jun 29 15:09:36 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291009.D Vial: 9
Acq On : 29 Jun 2010 11:40 Operator: MD
Sample : P1002127-026 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:09 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.07min 2280.549ng/ml m

response 326541913

6/30/10

6/30/10

49

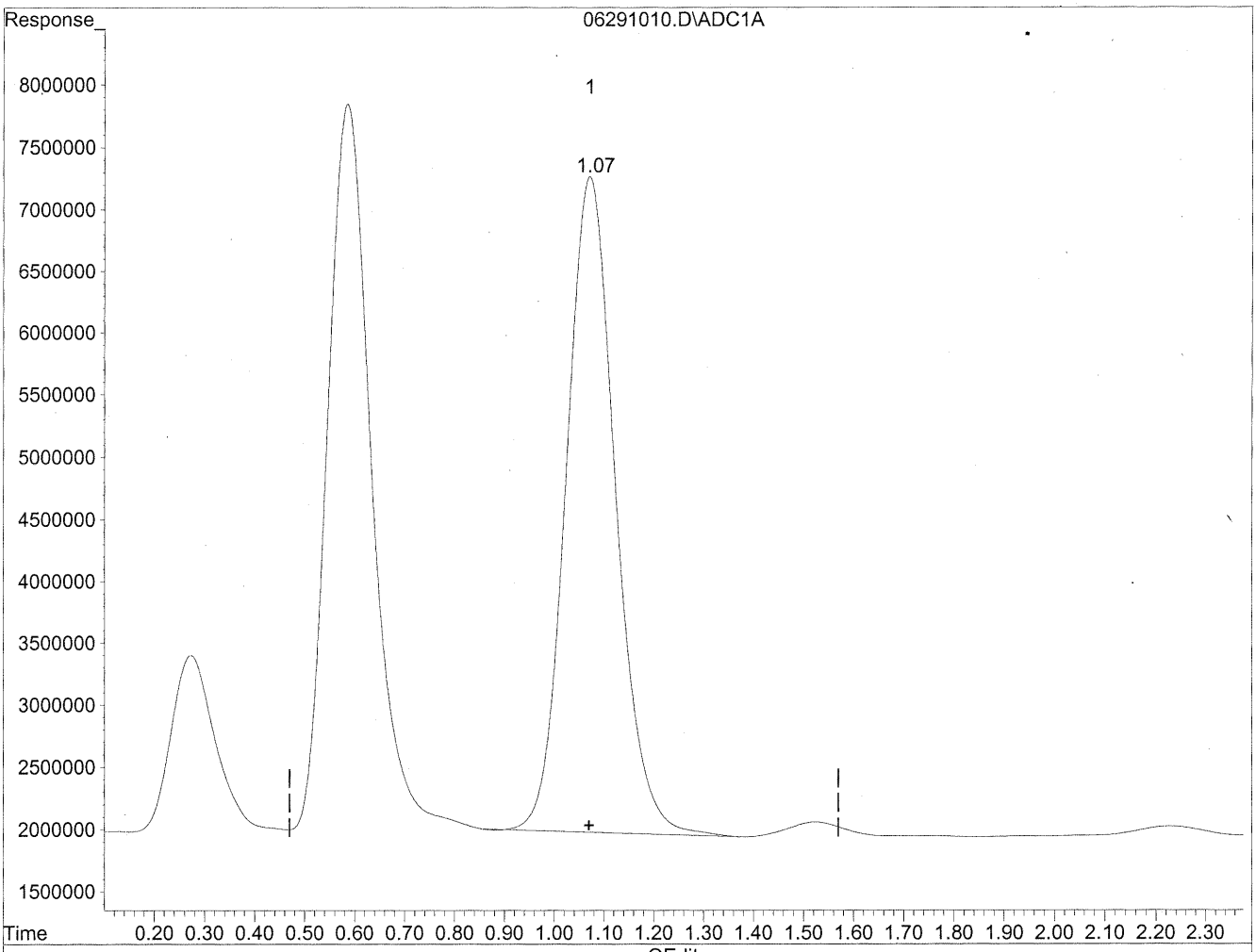
(+) = Expected Retention Time

06291009.D TO110610.M Tue Jun 29 15:09:40 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291010.D Vial: 10
Acq On : 29 Jun 2010 11:54 Operator: MD
Sample : P1002127-027 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:09 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.07min 2553.102ng/ml

response 365567643

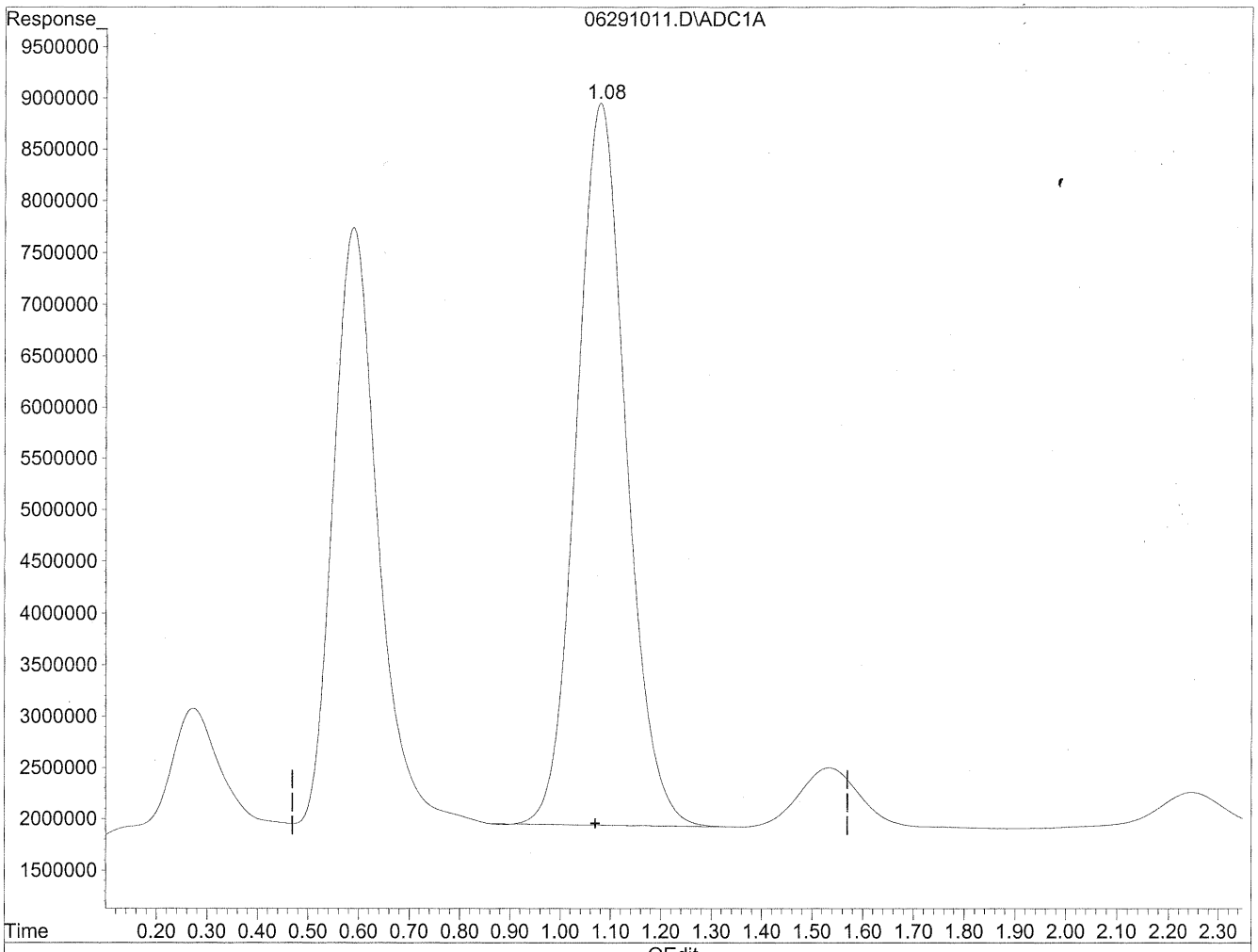
(+) = Expected Retention Time

06291010.D TO110610.M Tue Jun 29 15:09:57 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291011.D Vial: 11
Acq On : 29 Jun 2010 12:07 Operator: MD
Sample : P1002127-028 2.0ml 10x dil Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:10 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 3368.941ng/ml

response 482384082

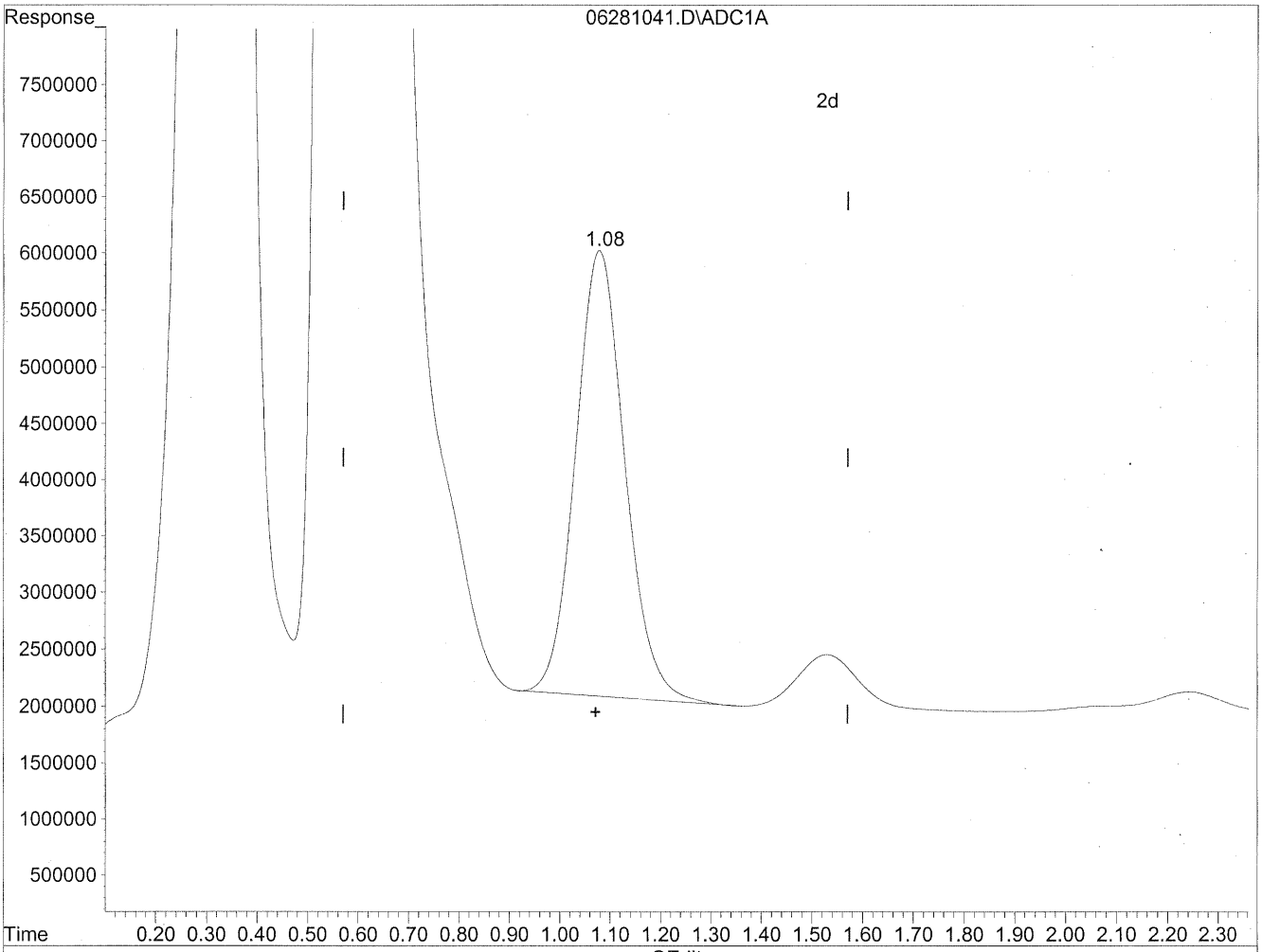
(+) = Expected Retention Time

06291011.D TO110610.M Tue Jun 29 15:10:17 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281041.D Vial: 41
Acq On : 28 Jun 2010 19:22 Operator: MD
Sample : P1002127-029 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 7:51 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



QEdit

(1) Formaldehyde

1.08min 1870.467ng/ml

response 267824086

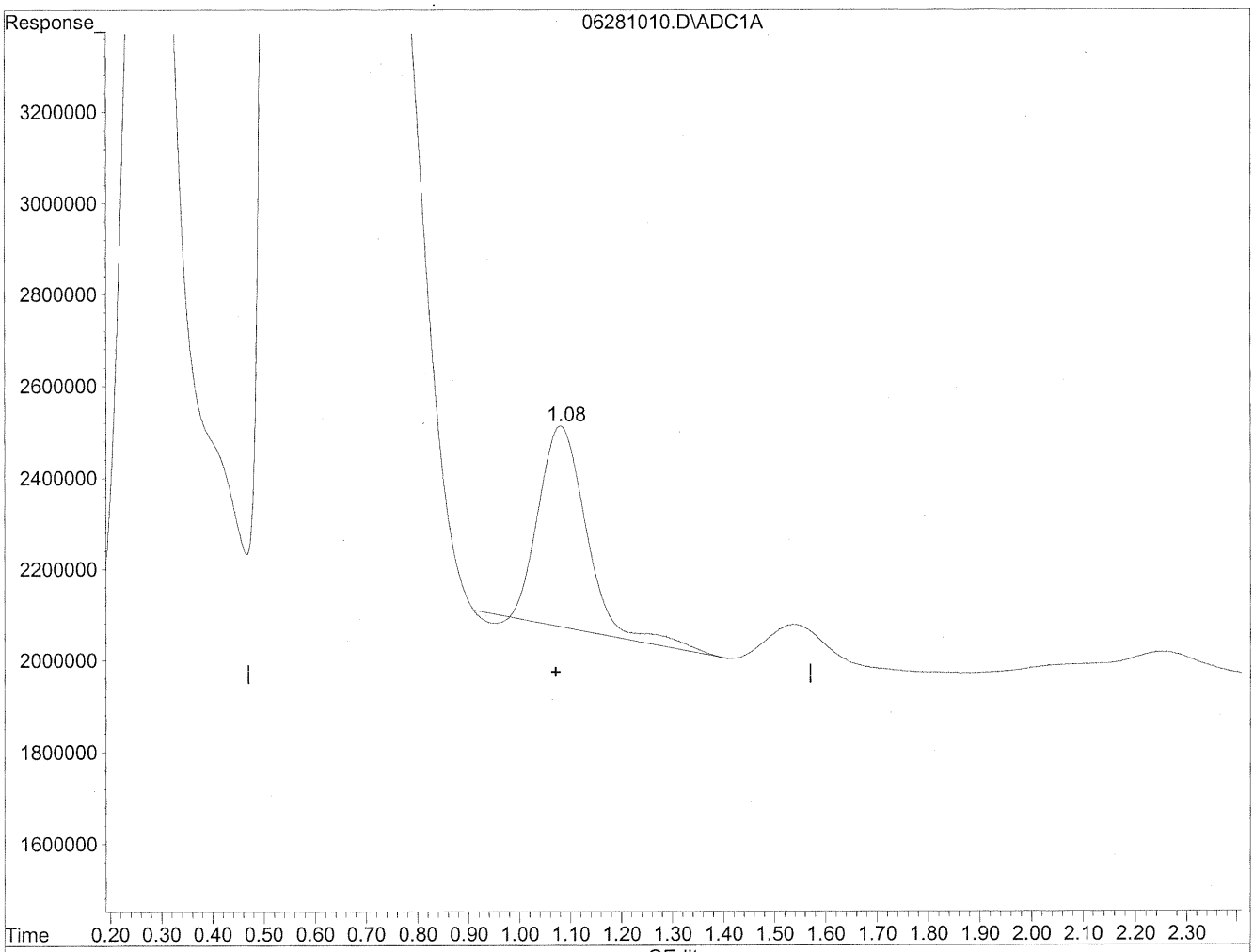
(+) = Expected Retention Time

06281041.D TO110610.M Tue Jun 29 13:27:46 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281010.D Vial: 10
Acq On : 28 Jun 2010 12:09 Operator: MD
Sample : P1002127-030 2.0ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 13:22 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 198.633ng/ml

response 28441368

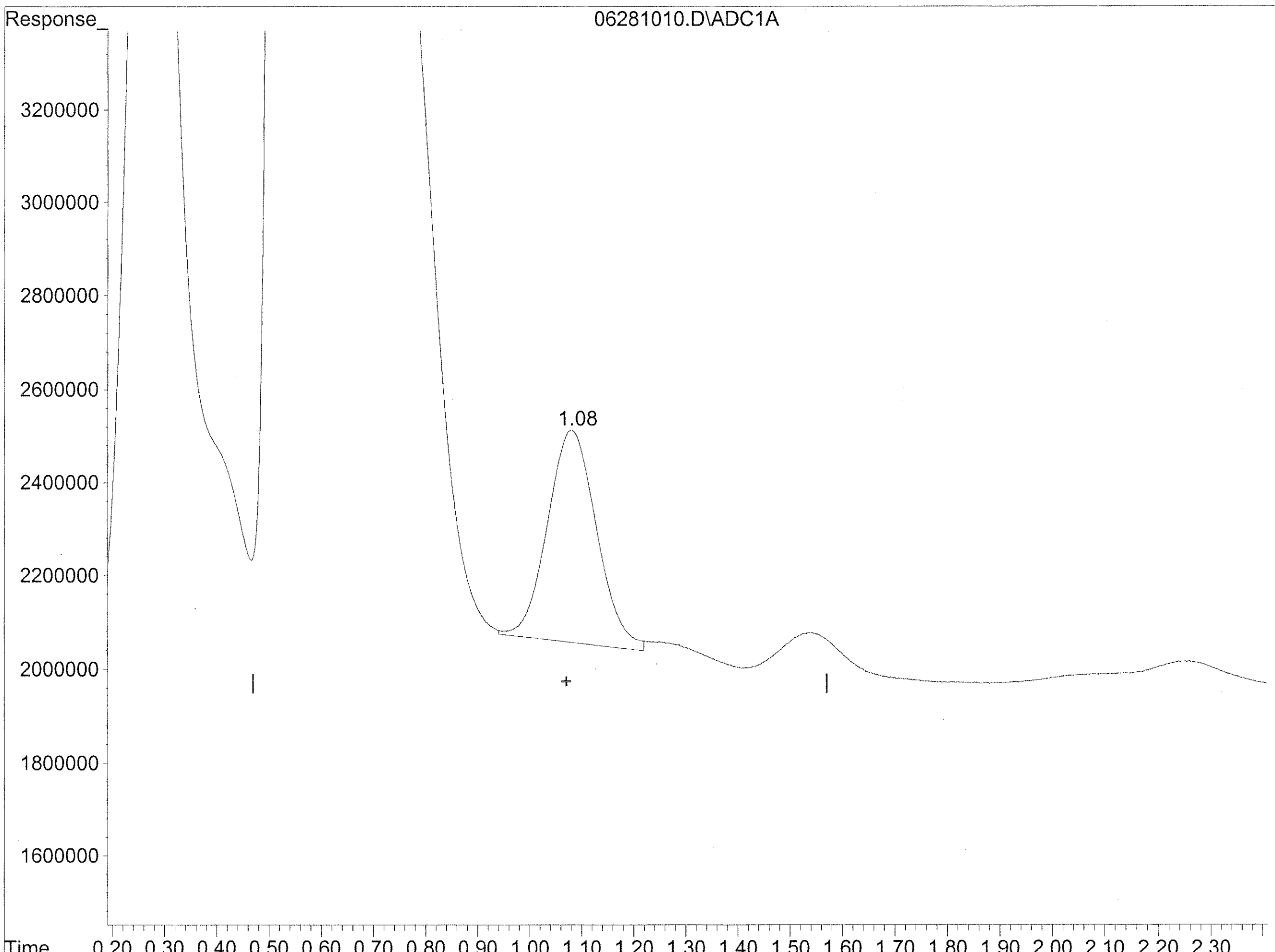
(+) = Expected Retention Time

06281010.D TO110610.M Tue Jun 29 13:22:17 2010

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2010_06\28\06281010.D      Vial: 10
Acq On    : 28 Jun 2010  12:09                       Operator: MD
Sample    : P1002127-030  2.0ml                      Inst    : LC 01
Misc      :                                          Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Jun 29 13:22 19110  Quant Results File: TO110610.RES

Method         : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title          : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update    : Tue Jun 29 12:58:25 2010
Response via   : Multiple Level Calibration
```



QEdit

(1) Formaldehyde

1.08min 209.147ng/ml m

response 29946921

ic 6/30/10

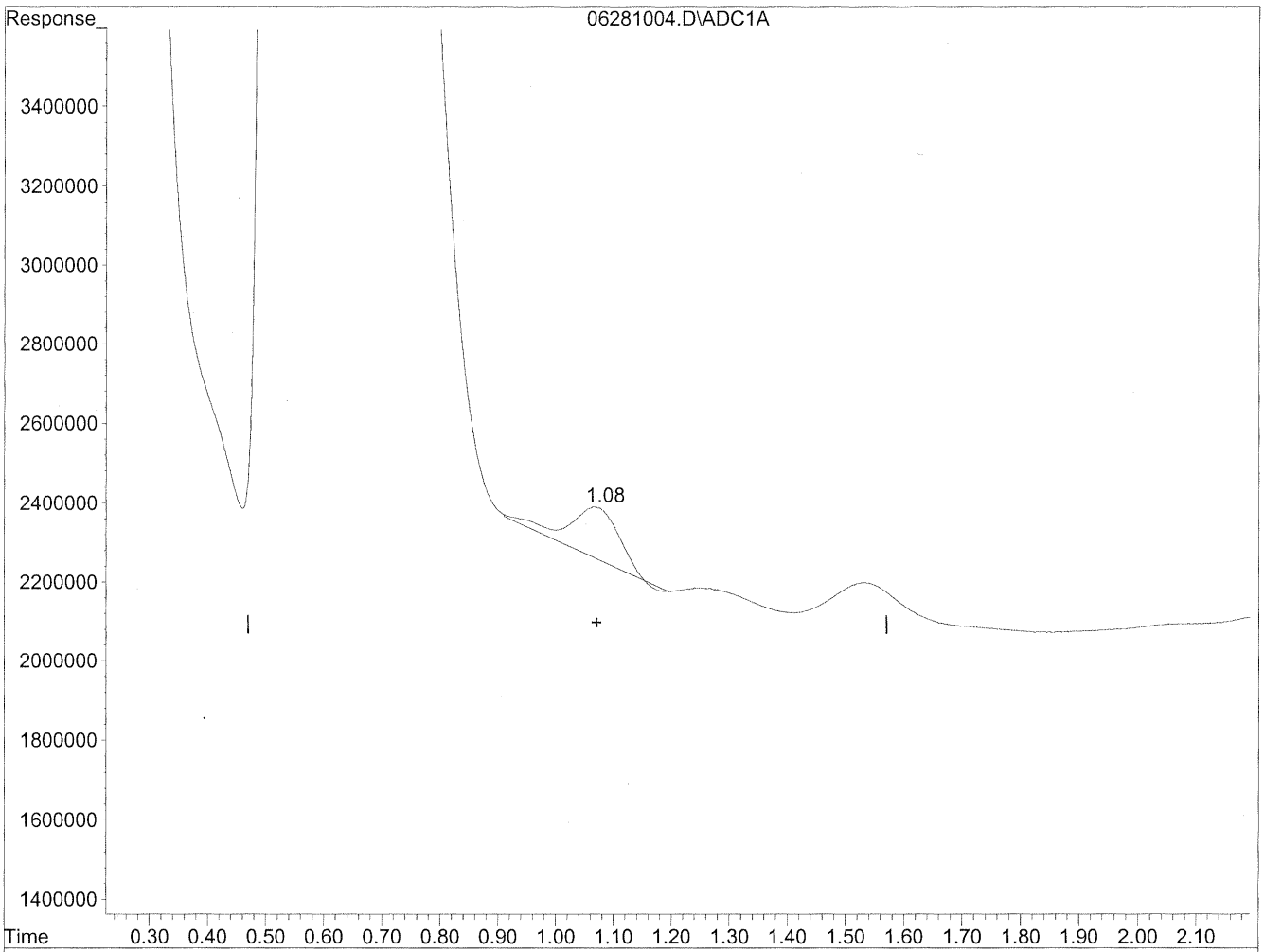
MD 6/30/10

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281004.D Vial: 4
Acq On : 28 Jun 2010 10:47 Operator: MD
Sample : MB 2.0ml Lot 09240 Inst : LC 01
Misc : Radiello passive sampler Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 12:59 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.07min 54.833ng/ml

response 7851368

(+) = Expected Retention Time

06281004.D TO110610.M Tue Jun 29 12:59:20 2010

INITIAL CALIBRATION STANDARDS

Response Factor Report LC 01

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Mar 21 12:19:47 2005

Calibration Files
50 =06241006.D 100 =06241009.D 500 =06241012.D
1500 =06241015.D 5000 =06241018.D 10 =06241021.D

	Compound	50	100	500	1500	5000	10	Avg		%RSD
1)	Formaldehyde	1.318	1.379	1.411	1.514	1.486	1.483	1.432	E5	5.25
2)	Acetaldehyde	1.058	1.081	1.079	1.153	1.124	1.121	1.103	E5	3.23
3)	Propionaldehyde	7.404	8.260	8.164	8.263	8.401	8.418	8.152	E4	4.65
4)	Crotonaldehyde	7.067	7.983	7.551	7.928	7.715	7.612	7.643	E4	4.31
5)	Butyraldehyde	6.916	7.832	7.126	6.824	7.197	7.117	7.169	E4	4.94
6)	Benzaldehyde	4.709	6.298	5.396	5.825	5.446	5.314	5.498	E4	9.68
7)	Isovaleraldehyde	5.460	7.207	6.289	6.728	6.364	6.250	6.383	E4	9.07
8)	Valeraldehyde	5.513	6.252	5.544	5.530	5.665	5.685	5.698	E4	4.93
9)	o-Tolualdehyde	4.516	4.892	4.810	5.078	4.875	4.804	4.829	E4	3.79
10)	m,p-Tolualdehyde	4.137	4.321	4.226	4.505	4.383	4.337	4.318	E4	2.95
11)	Hexaldehyde	4.787	5.800	5.158	5.140	5.374	5.281	5.257	E4	6.33
12)	2,5-Dimethylbenzald	3.676	4.598	3.679	3.946	3.822	3.766	3.915	E4	8.93

#) = Out of Range

COLUMBIA ANALYTICAL SERVICES, INC.

Method: TO-11A
Analyst: MD

Printed : 06/25/10
Instrument : LC#1
Date Analysis : 06/24/10
Detector : UV-VIS 360
Sample Amount : 5ul

6/25/10 +IL 6/28/10

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 TO-11A S2	6762023	2.57%	5145025	2.75%	3983499	7.61%	3778559	6.93%	3523246	1.89%	2365205	0.45%
50ng/ml TO-11A S2	6598476	0.09%	5336318	0.86%	3364453	9.11%	3240110	8.31%	3652578	5.63%	2527622	7.35%
50ng/ml TO-11A S2	6416946	2.66%	5390444	1.89%	3757337	1.50%	3582407	1.38%	3198175	7.51%	2170695	7.81%
100ng/ml TO-11A S	13684484	0.80%	10561892	2.30%	8914731	7.93%	7957674	0.32%	7348500	6.17%	6214018	1.34%
100ng/ml TO-11A S	13511387	2.05%	10740786	0.64%	8027964	2.80%	7500614	6.04%	8049959	2.79%	6787382	7.76%
100ng/ml TO-11A S	14186865	2.85%	11127714	2.94%	7835935	5.13%	8490383	6.36%	8096555	3.38%	5893577	6.43%
500ng/ml TO-11A S	71248470	1.02%	54504441	1.07%	41248067	1.05%	38090464	0.89%	35410354	0.62%	28344164	5.05%
500ng/ml TO-11A S	70508340	0.03%	54114192	0.34%	40426693	0.96%	38548382	2.11%	35949330	0.89%	26799052	0.67%
500ng/ml TO-11A S	69824657	1.00%	53171079	1.41%	40786533	0.08%	36622016	3.00%	35535481	0.27%	25798064	4.38%
1500ng/ml TO-11A	230593016	1.56%	175356392	1.41%	125879750	1.56%	120696891	1.50%	103399084	1.02%	89107201	1.99%
1500ng/ml TO-11A	225421129	0.72%	172419255	0.29%	124283598	0.27%	119167899	0.21%	101934633	0.41%	86843740	0.60%
1500ng/ml TO-11A	225129784	0.84%	170977797	1.12%	121684007	1.83%	116879696	1.71%	101737279	0.61%	86160446	1.38%
5000ng/ml TO-11A	752167389	1.22%	565379435	0.58%	421481275	0.34%	387209880	0.37%	360334815	0.14%	271843062	0.17%
5000ng/ml TO-11A	741687155	0.19%	560621731	0.26%	418074248	0.47%	384212117	0.40%	359182440	0.18%	271646285	0.24%
5000ng/ml TO-11A	735352713	1.04%	560284478	0.32%	420570615	0.13%	385901475	0.03%	359960551	0.04%	273410909	0.41%
10000ng/ml TO-11A	1506388672	1.59%	1143475045	2.03%	857670773	1.88%	774502971	1.75%	723112269	1.60%	541585412	1.92%
10000ng/ml TO-11A	1465408979	1.18%	1108342081	1.11%	834978009	0.81%	755899749	0.70%	707762167	0.56%	527522590	0.73%
10000ng/ml TO-11A	1476855713	0.41%	1110424466	0.92%	832841780	1.07%	753218148	1.05%	704333220	1.04%	525054783	1.19%

COLUMBIA ANALYTICAL SERVICES, INC.

Method: TO-11A
Analyst: MD

Printed : 06/25/10
Instrument : LC#1
Date Analysis : 06/24/10
Detector : UV-VIS 360
Sample Amount : 5ul

CALIBRATION RESPONSE FACTOR SUMMARY

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 TO-11A S2	2265328	17.02%	2716276	1.46%	2373630	5.12%	4098785	0.91%	2290365	4.32%	1619465	11.90%
50ng/ml TO-11A S2	2995744	9.74%	2768346	0.43%	2334081	3.37%	3866830	6.52%	2424895	1.30%	2173046	18.22%
50ng/ml TO-11A S2	2928376	7.27%	2784817	1.03%	2066036	8.50%	4443968	7.43%	2465808	3.01%	1722129	6.32%
100ng/ml TO-11A S	7015814	2.65%	5776234	7.60%	5309971	8.54%	9000241	4.15%	5629872	2.94%	4497151	2.20%
100ng/ml TO-11A S	7873759	9.26%	6332528	1.29%	4800335	1.88%	8786398	1.67%	5891760	1.58%	4437325	3.50%
100ng/ml TO-11A S	6729977	6.61%	6646232	6.31%	4566176	6.66%	8138944	5.82%	5878688	1.35%	4860735	5.70%
500ng/ml TO-11A S	32418035	3.09%	28548123	2.98%	24943255	3.71%	43217145	2.27%	27064970	4.95%	18911052	2.80%
500ng/ml TO-11A S	31444882	0.00%	28285881	2.03%	23916376	0.56%	42173453	0.20%	25614342	0.67%	18988988	3.23%
500ng/ml TO-11A S	30475379	3.09%	26332477	5.01%	23290726	3.16%	41386366	2.06%	24685885	4.28%	17286981	6.03%
1500ng/ml TO-11A	102723576	1.78%	84987047	2.46%	77943499	2.33%	138317069	2.35%	79547312	3.18%	60933239	2.93%
1500ng/ml TO-11A	100332819	0.58%	81906561	1.26%	73725386	3.21%	134049332	0.81%	75378424	2.22%	58178943	1.72%
1500ng/ml TO-11A	99712150	1.20%	81955851	1.20%	76844694	0.88%	133062119	1.54%	76355051	0.96%	58479212	1.21%
5000ng/ml TO-11A	318898458	0.22%	283824088	0.21%	245490595	0.72%	439534493	0.28%	269904263	0.44%	191751893	0.34%
5000ng/ml TO-11A	316239705	0.62%	281768408	0.52%	241706915	0.84%	435865650	0.56%	266567105	0.80%	189585119	0.80%
5000ng/ml TO-11A	319488895	0.40%	284129216	0.31%	244039778	0.12%	439548558	0.28%	269656485	0.35%	191978970	0.46%
10000ng/ml TO-11A	637094049	1.93%	579085929	1.87%	490639974	2.13%	882689956	1.76%	537444487	1.77%	383332870	1.78%
10000ng/ml TO-11A	619285794	0.92%	564056973	0.78%	475987748	0.92%	860464776	0.80%	525177534	0.55%	373842408	0.74%
10000ng/ml TO-11A	618766885	1.01%	562282799	1.09%	474574515	1.21%	859080340	0.96%	521619735	1.22%	372739681	1.04%

AVERAGE RESPONSE FACTOR

	Form-Aldehyde	Acet-Aldehyde	Propion-Aldehyde	Croton-Aldehyde	Butyr-Aldehyde	Benz-Aldehyde
50ng/ml TO-11A S	6592482	5290596	3701763	3533692	3458000	2354507
100ng/ml TO-11A S	13794245	10810131	8259543	7982890	7831671	6298326
500ng/ml TO-11A S	70527156	53929904	40820431	37753621	35631722	26980427
1500ng/ml TO-11A	227047976	172917815	123949118	118914829	102356999	87370462
5000ng/ml TO-11A	743069086	562095215	420042046	385774491	359825935	272300085
10000ng/ml TO-11	1482884455	1120747197	841830187	761206956	711735885	531387595

AVERAGE RESPONSE FACTOR

	Isovaler-Aldehyde	Valer-Aldehyde	o-Tolu-Aldehyde	m,p-Tolu-Aldehyde	Hex-Aldehyde	2,5-Dimethl benz-Aldehyde
50ng/ml TO-11A S	2729816	2756480	2257916	4136528	2393689	1838213
100ng/ml TO-11A S	7206517	6251665	4892161	8641861	5800107	4598404
500ng/ml TO-11A S	31446099	27722160	24050119	42258988	25788399	18395674
1500ng/ml TO-11A	100922848	82949820	76171193	135142840	77093596	59197131
5000ng/ml TO-11A	318209019	283240571	243745763	438316234	268709284	191105327
10000ng/ml TO-11	625048909	568475234	480400746	867411691	528080585	376638320

COMPOUND	50	100	500	1500	5000	10000	AVERAGE	SD	%RSD
Formaldehyde	131849.6333	1.38E+05	1.41E+05	1.51E+05	1.49E+05	1.48E+05	1.43E+05	7.52E+03	5.25%
Acetaldehyde	105811.9133	1.08E+05	1.08E+05	1.15E+05	1.12E+05	1.12E+05	1.10E+05	3.56E+03	3.23%
Propionaldehyde	74035.26	8.26E+04	8.16E+04	8.26E+04	8.40E+04	8.42E+04	8.15E+04	3.79E+03	4.65%
Crotonaldehyde	70673.84	7.98E+04	7.55E+04	7.93E+04	7.72E+04	7.61E+04	7.64E+04	3.29E+03	4.31%
Butyraldehyde	69159.99333	7.83E+04	7.13E+04	6.82E+04	7.20E+04	7.12E+04	7.17E+04	3.54E+03	4.94%
Benzaldehyde	47090.14667	6.30E+04	5.40E+04	5.82E+04	5.45E+04	5.31E+04	5.50E+04	5.32E+03	9.68%
Isovaleraldehyde	54596.32	7.21E+04	6.29E+04	6.73E+04	6.36E+04	6.25E+04	6.38E+04	5.79E+03	9.07%
Valeraldehyde	55129.59333	6.25E+04	5.54E+04	5.53E+04	5.66E+04	5.68E+04	5.70E+04	2.81E+03	4.93%
o-Tolualdehyde	45158.31333	4.89E+04	4.81E+04	5.08E+04	4.87E+04	4.80E+04	4.83E+04	1.83E+03	3.79%
m,p-Tolualdehyde	41365.27667	4.32E+04	4.23E+04	4.50E+04	4.38E+04	4.34E+04	4.32E+04	1.27E+03	2.95%
Hexaldehyde	47873.78667	5.80E+04	5.16E+04	5.14E+04	5.37E+04	5.28E+04	5.26E+04	3.33E+03	6.33%
2,5-Dimethylbenzaldehyde	36764.26667	4.60E+04	3.68E+04	3.95E+04	3.82E+04	3.77E+04	3.91E+04	3.50E+03	8.93%

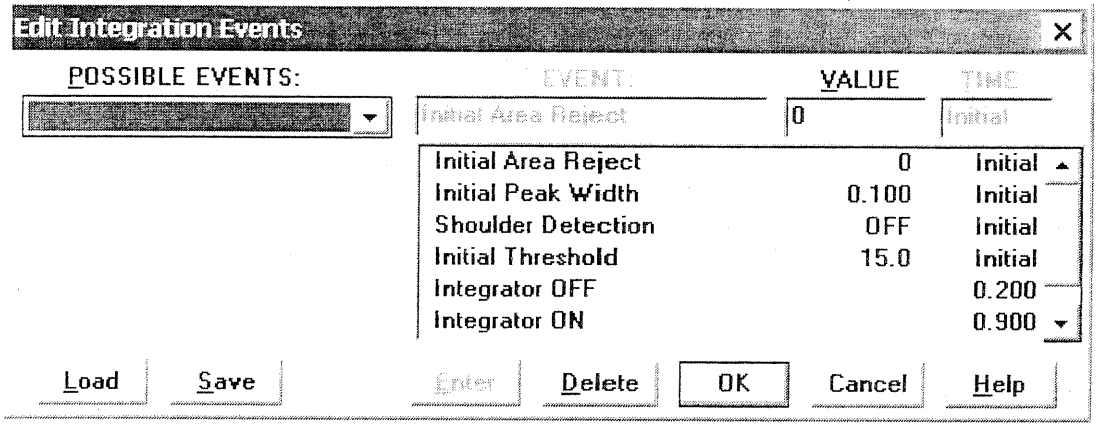
TO-11A CALIBRATION STANDARDS LIST
50ng/ml TO-11A S21-03091012
100ng/ml TO-11A S21-03091009
500ng/ml TO-11A S21-03091008
1500ng/ml TO-11A S21-04211003
5000ng/ml TO-11A S21-03091001
10000ng/ml TO-11A S21-03091007

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Mon Mar 21 12:19:47 2005
Response via : Initial Calibration

#	ID	Conc	ISTD Conc	Path\File
1	50	50.00	0.00	J:\LC01\DATA\TO11\2010_06\24\06241006.D
2	100	100.00	0.00	J:\LC01\DATA\TO11\2010_06\24\06241009.D
3	500	500.00	0.00	J:\LC01\DATA\TO11\2010_06\24\06241012.D
4	1500	1500.00	0.00	J:\LC01\DATA\TO11\2010_06\24\06241015.D
5	5000	5000.00	0.00	J:\LC01\DATA\TO11\2010_06\24\06241018.D
6	10	10000.00	0.00	J:\LC01\DATA\TO11\2010_06\24\06241021.D

#	ID	Update Time	Quant Time	Acquisition Time
1	50	Jun 24 16:53 2010	Jun 24 13:35 19110	24 Jun 2010 12:31
2	100	Jun 24 16:53 2010	Jun 24 16:47 19110	24 Jun 2010 13:12
3	500	Jun 24 16:53 2010	Jun 24 14:27 19110	24 Jun 2010 13:52
4	1500	Jun 24 16:54 2010	Jun 24 15:18 19110	24 Jun 2010 14:33
5	5000	Jun 24 16:54 2010	Jun 24 15:29 19110	24 Jun 2010 15:13
6	10	Jun 24 16:55 2010	Jun 24 16:08 19110	24 Jun 2010 15:54

TO110610.M Fri Jun 25 09:13:40 2010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EVENT:
VALUE
TIME:
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Initial
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Initial Threshold 15.0 Initial
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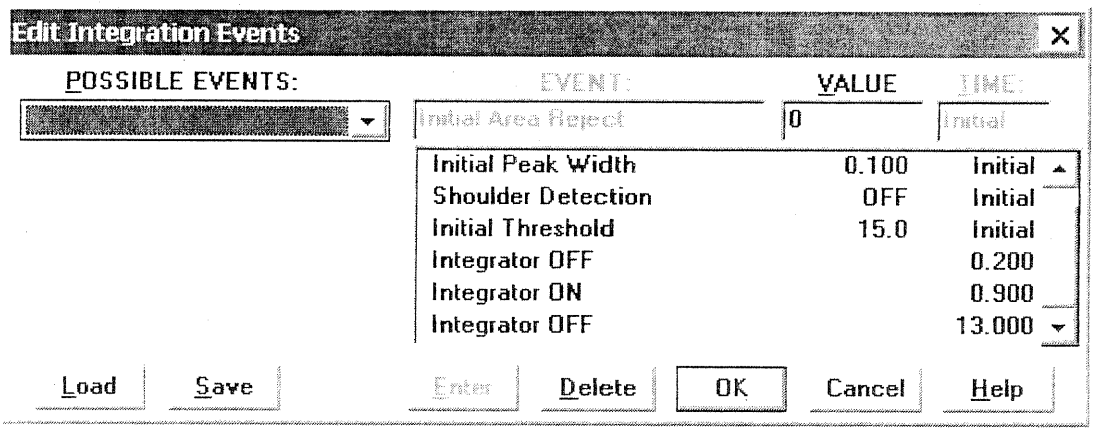
Integrator ON 0.900
Load
Save
Enter
Delete
OK
Cancel
Help


Edit Integration Events
POSSIBLE EVENTS:
EVENT:
VALUE
TIME:
Initial Area Reject
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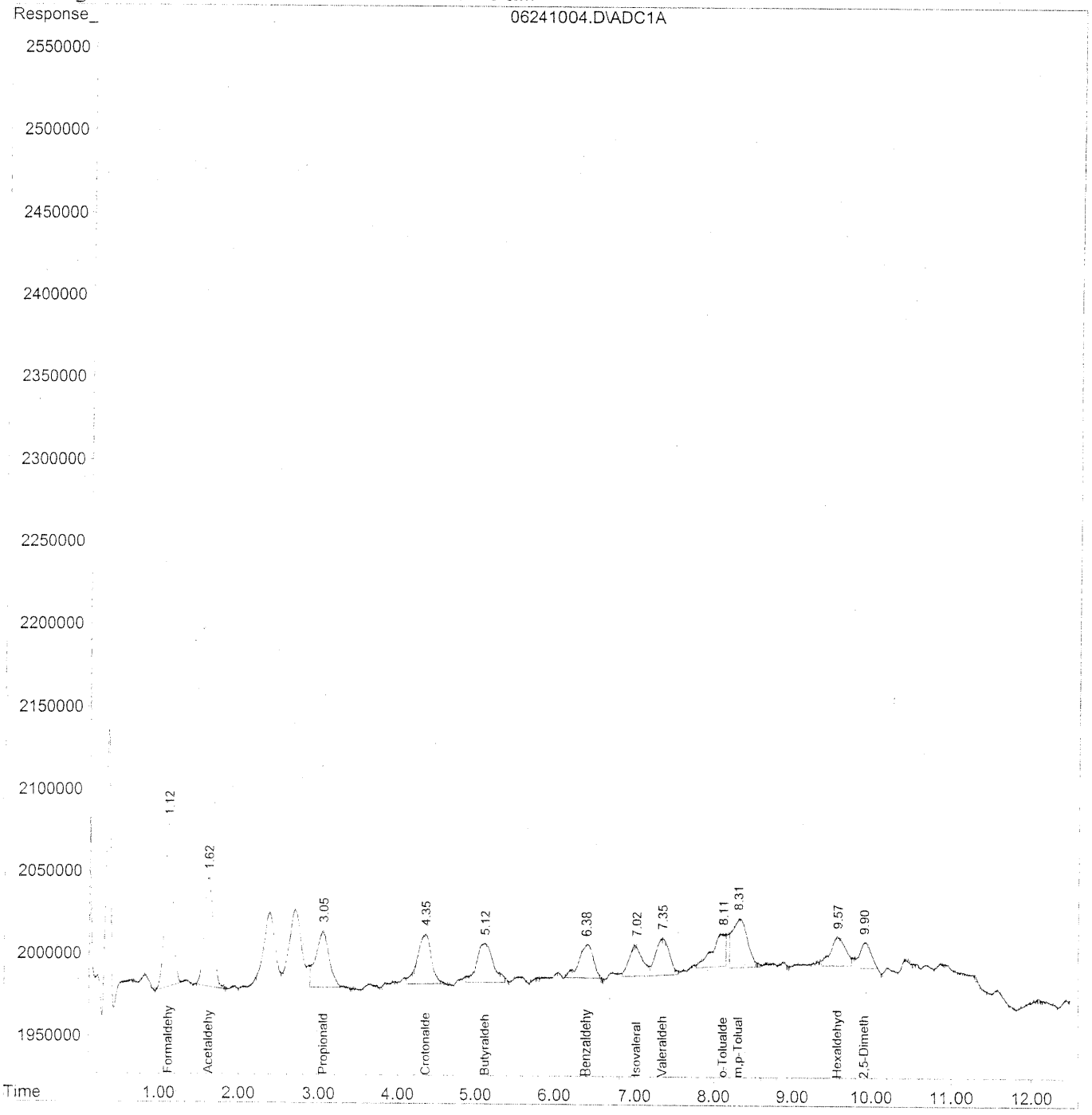
Initial
Initial Peak Width 0.100 Initial
Shoulder Detection OFF Initial
Initial Threshold 15.0 Initial
Integrator OFF 0.200
Integrator ON 0.900
Integrator OFF 13.000
Load
Save
Enter
Delete
OK
Cancel
Help

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:41 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

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



Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:41 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Initial Calibration
DataAcq Meth : TO11S.M

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

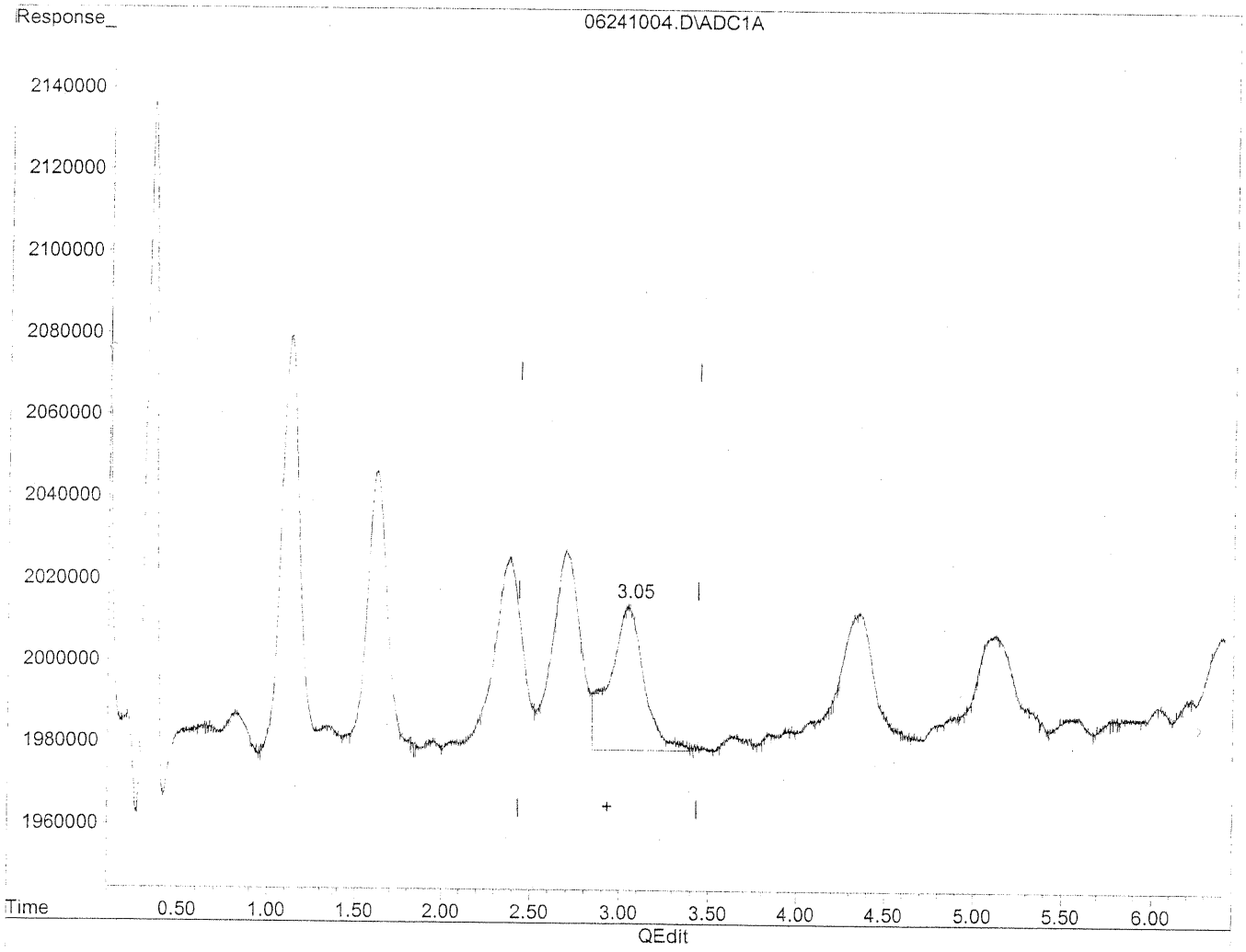
	Compound	R.T.	Response	Conc Units
	Target Compounds			
1)	Formaldehyde	1.12	6762023	30.366 ng/ml
2)	Acetaldehyde	1.62	5145025	30.190 ng/ml
3)	Propionaldehyde	3.05	3983499	30.878 ng/mlm
4)	Crotonaldehyde	4.35	3778559	32.240 ng/mlm
5)	Butyraldehyde	5.12	3523246	32.448 ng/mlm
6)	Benzaldehyde	6.38	2365205	30.376 ng/mlm
7)	Isovaleraldehyde	7.02	2265328	24.701 ng/mlm
8)	Valeraldehyde	7.35	2716276	31.385 ng/mlm
9)	o-Tolualdehyde	8.11	2373630	33.301 ng/mlm
10)	m,p-Tolualdehyde	8.31	4098785	64.001 ng/mlm
11)	Hexaldehyde	9.57f	2290365	27.309 ng/mlm
12)	2,5-Dimethylbenzaldehyde	9.90	1619465	28.403 ng/mlm

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(3) Propionaldehyde

3.04min 37.475ng/ml

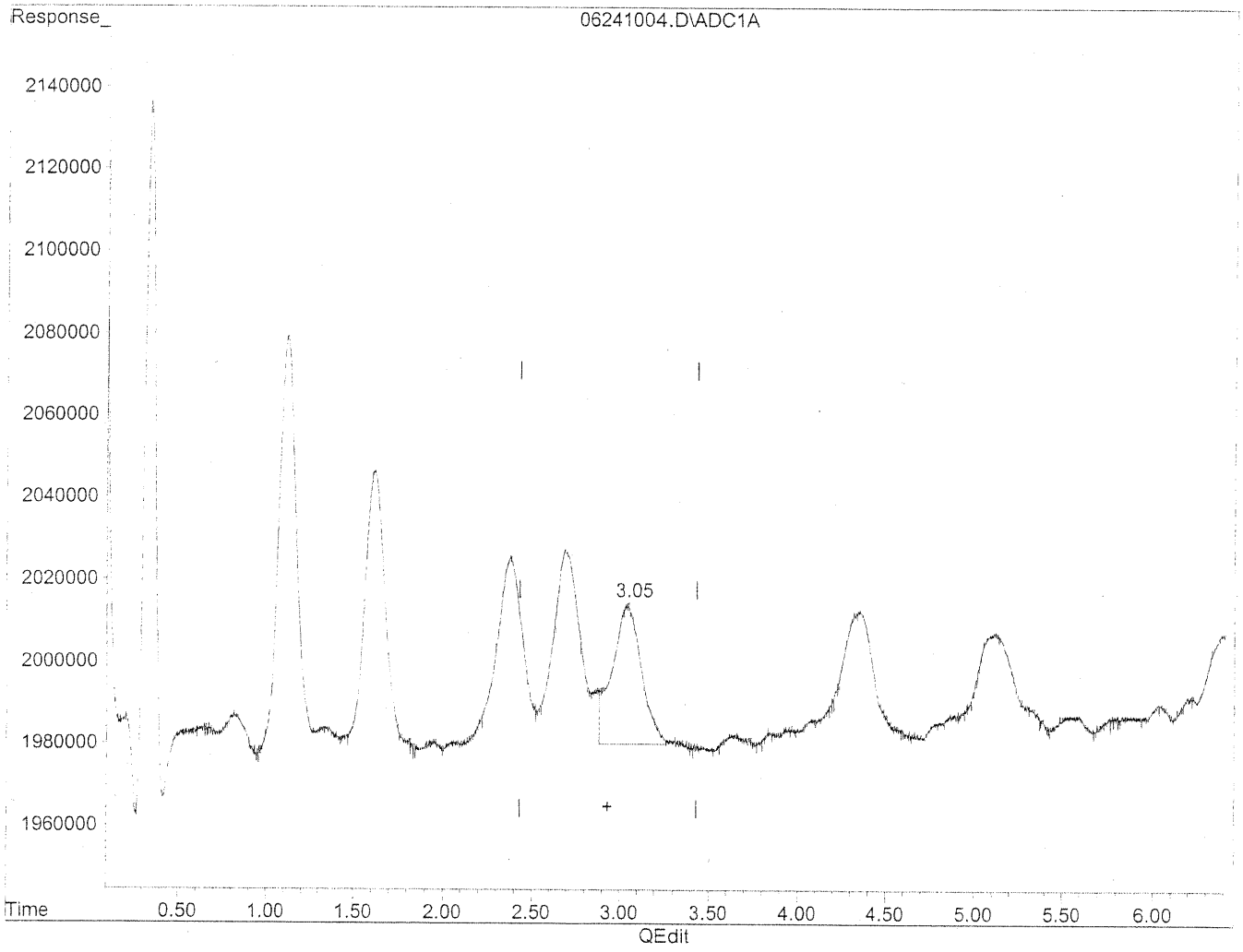
response 4834674

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(3) Propionaldehyde

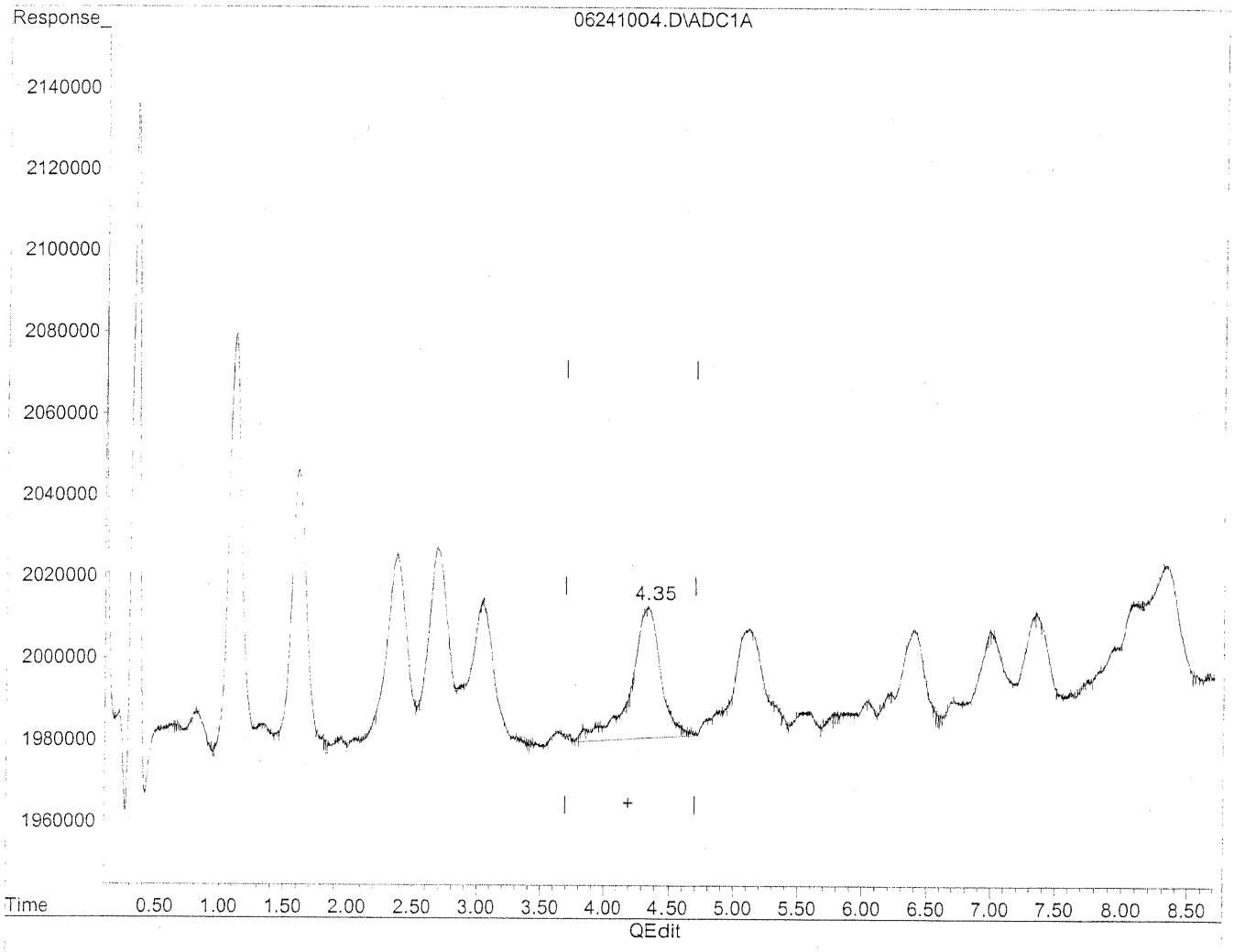
3.05min 30.878ng/ml m

response 3983499

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.35min 43.068ng/ml

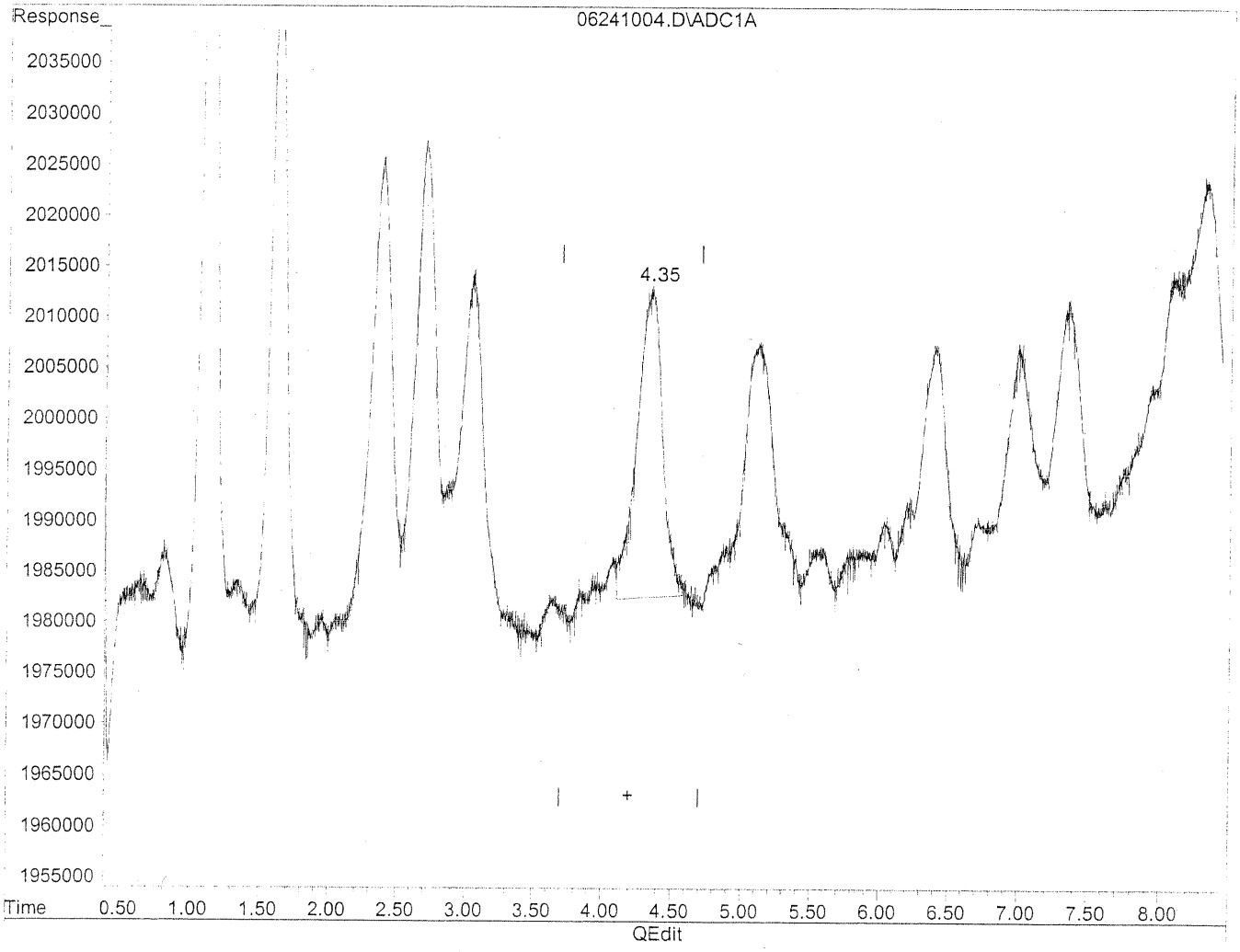
response 5047623

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.35min 32.240ng/ml m

response 3778559

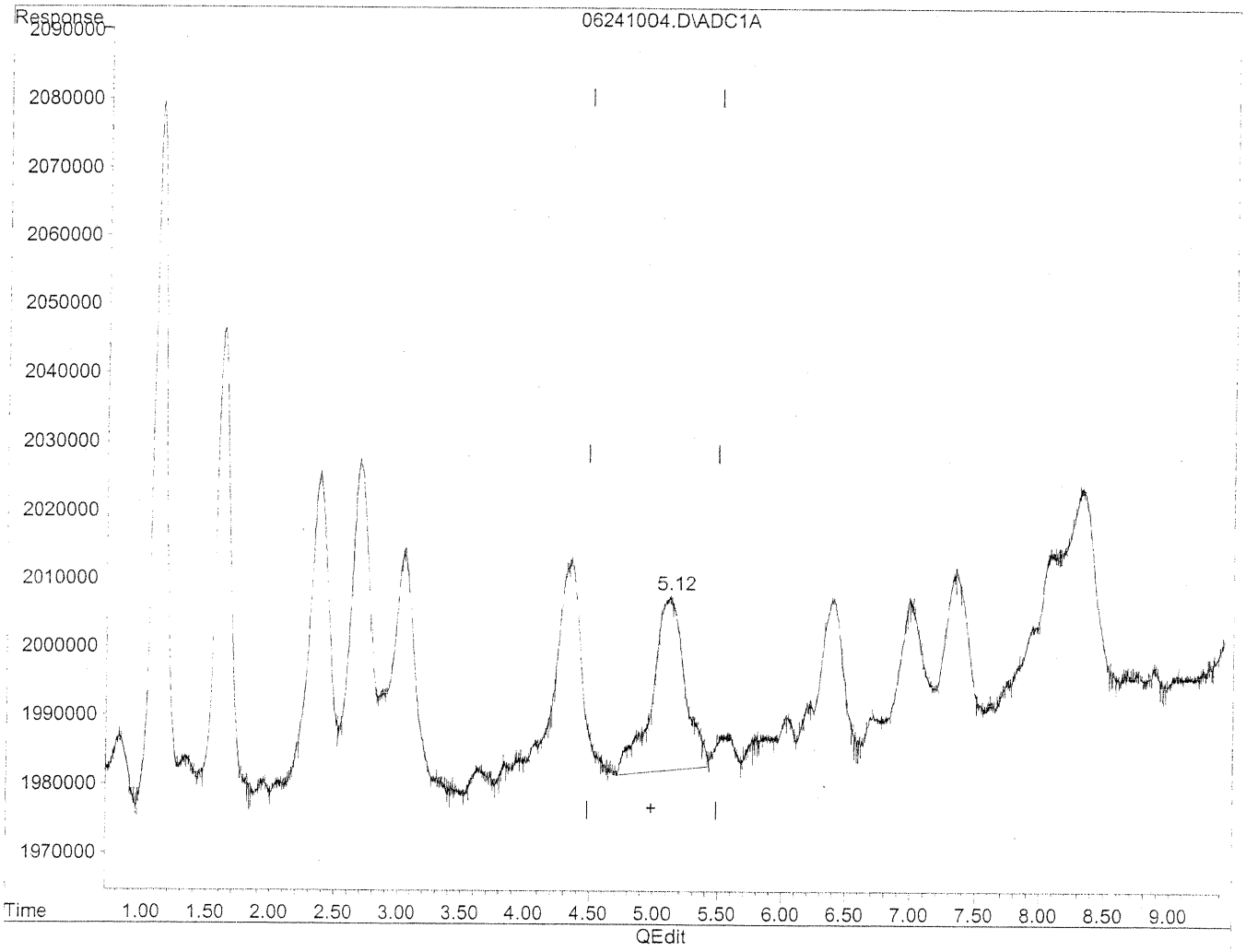
TC 6/28/10

TC MD 6/25/10

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.12min 40.066ng/ml

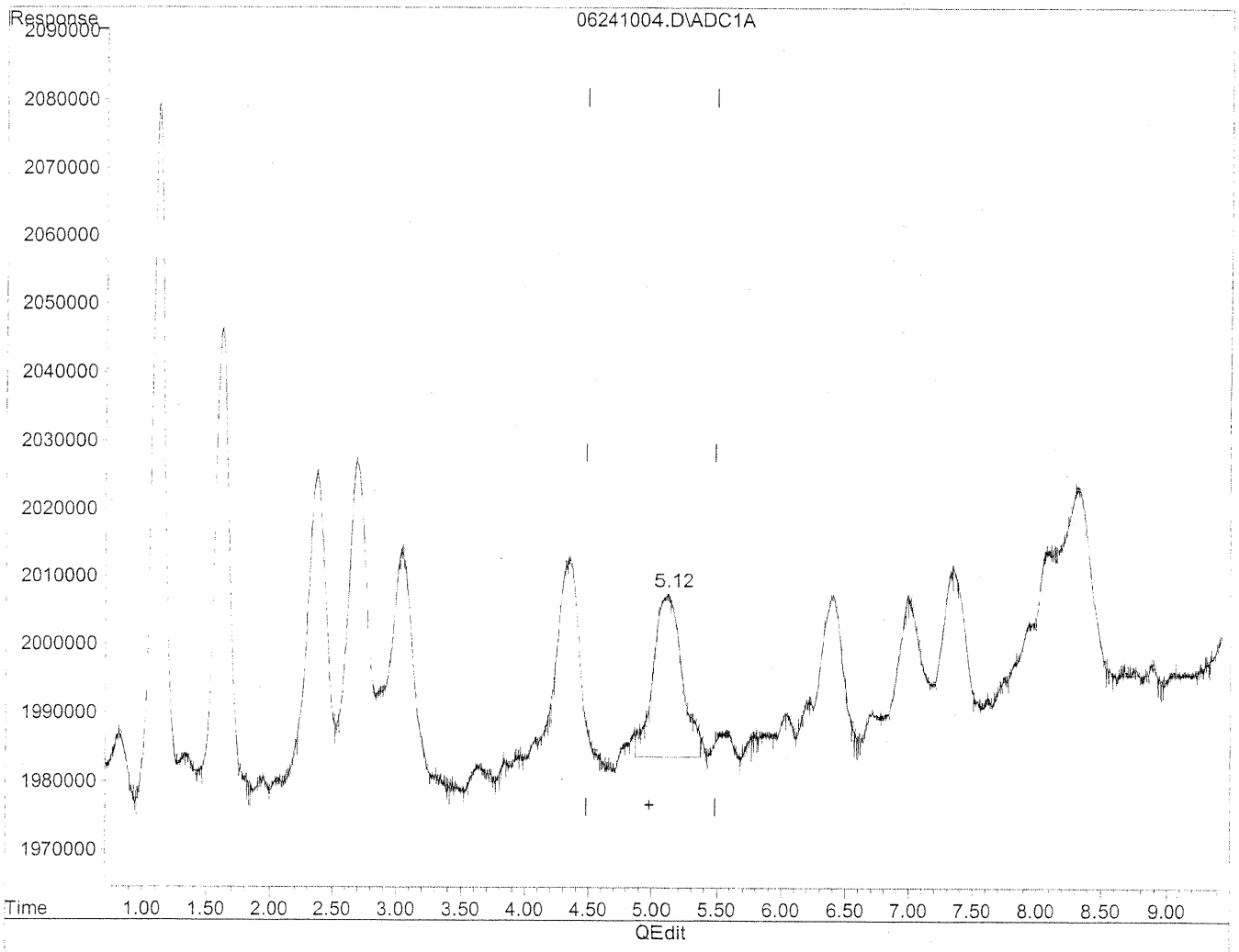
response 4350362

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(5) Butyraldehyde

5.12min 32.448ng/ml m

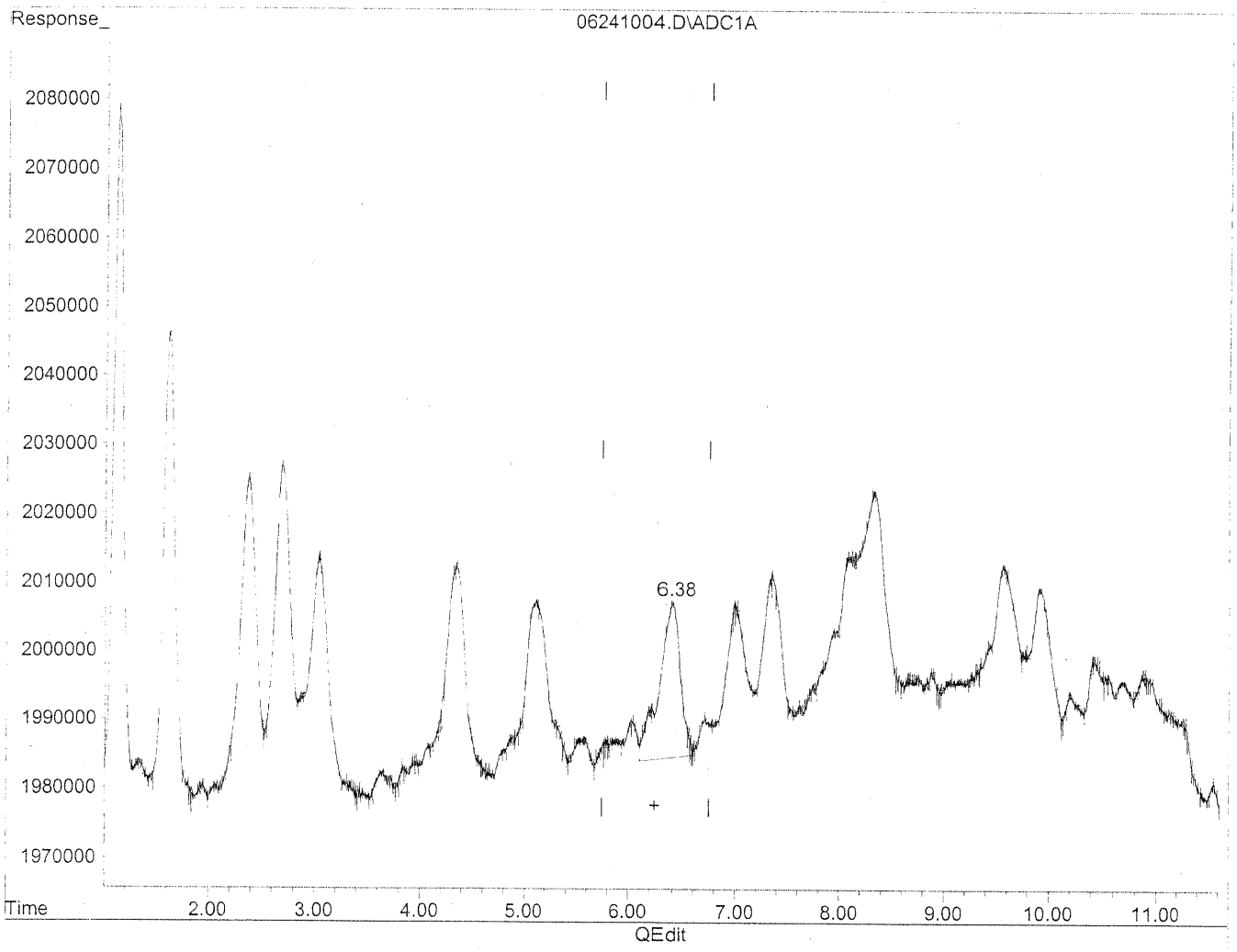
response 3523246

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.40min 39.370ng/ml

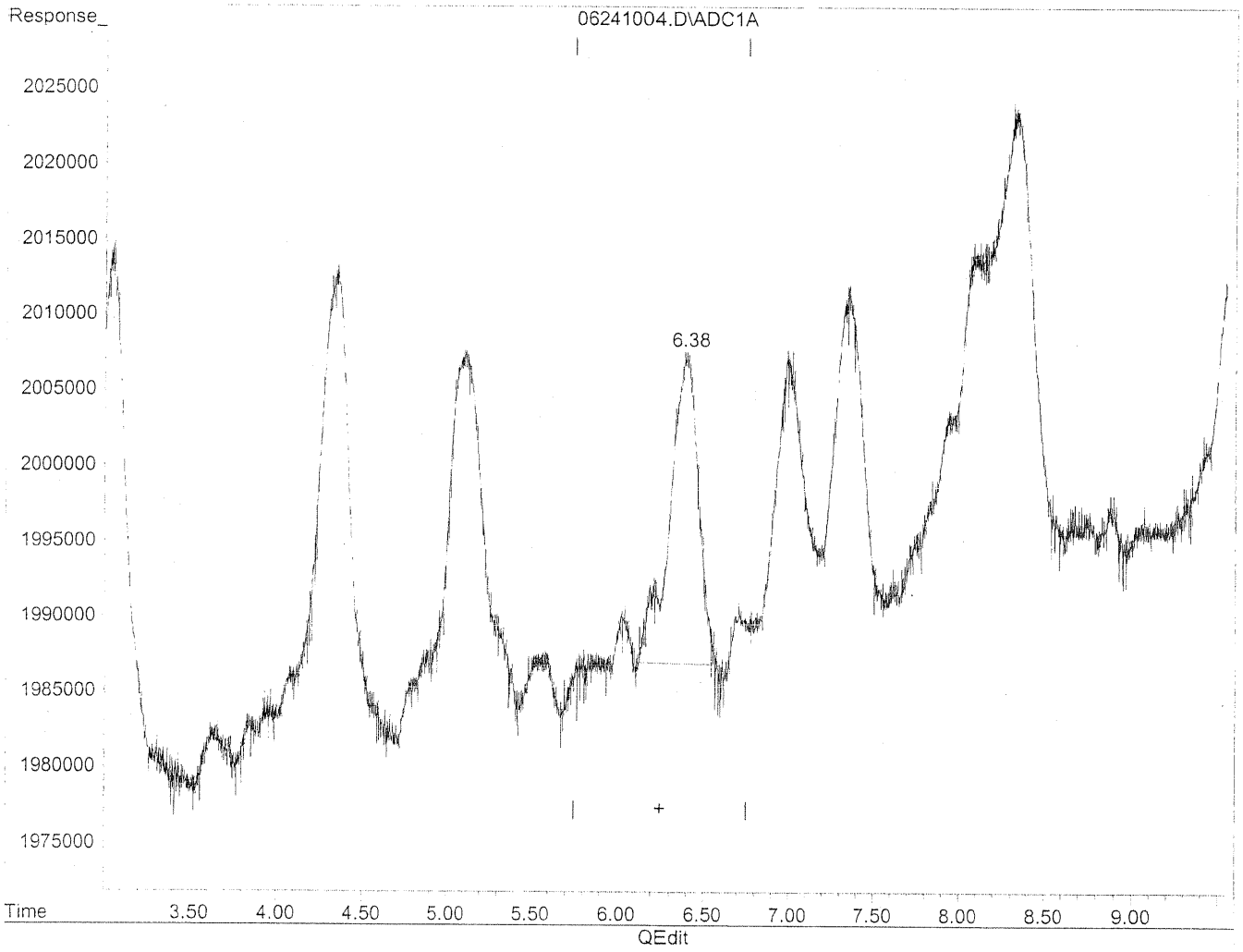
response 3065495

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 16:32:35 2010
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.38min 30.376ng/ml m

response 2365205

HC 6/28/10

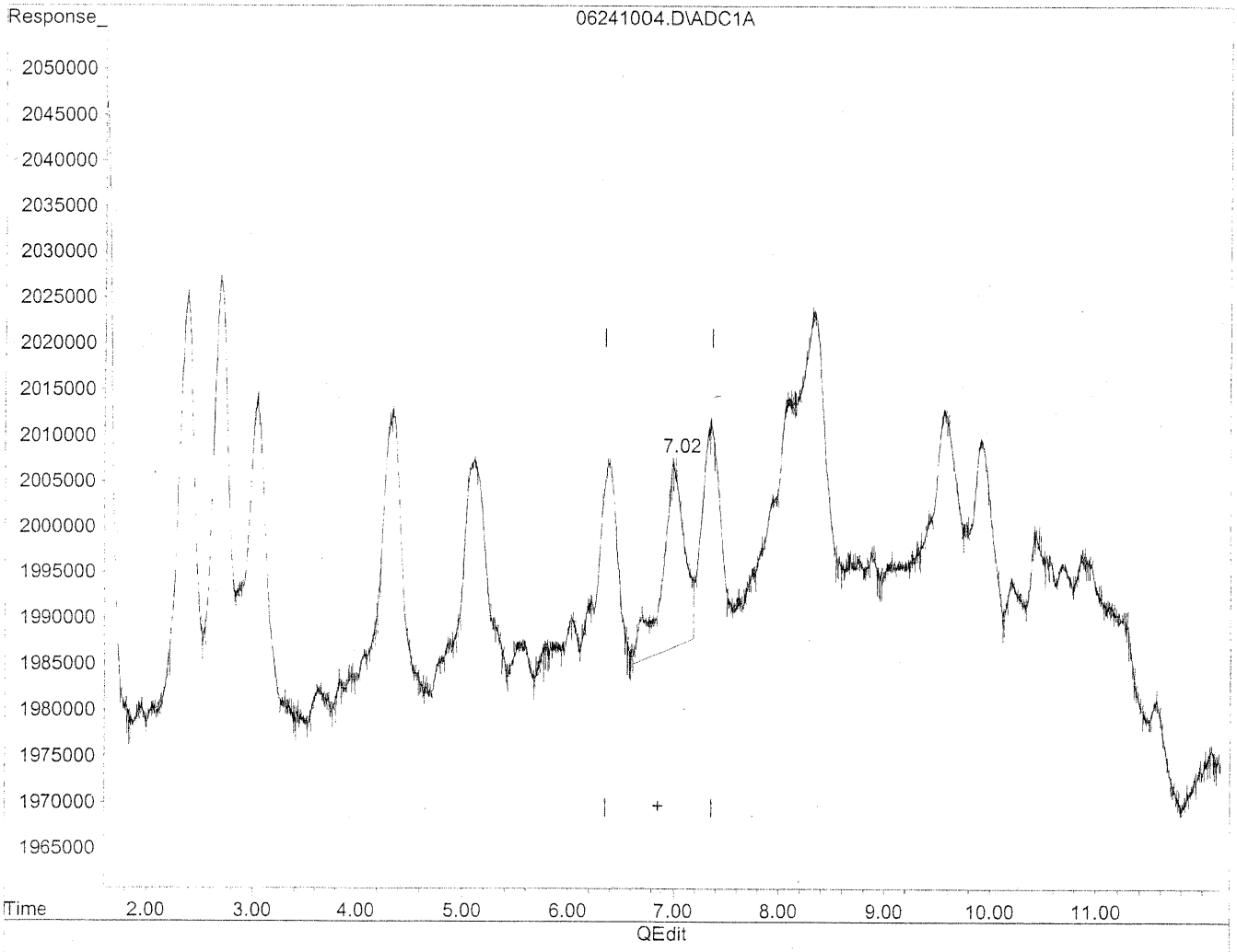
JZ 6/28/10

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

6.99min 32.048ng/ml

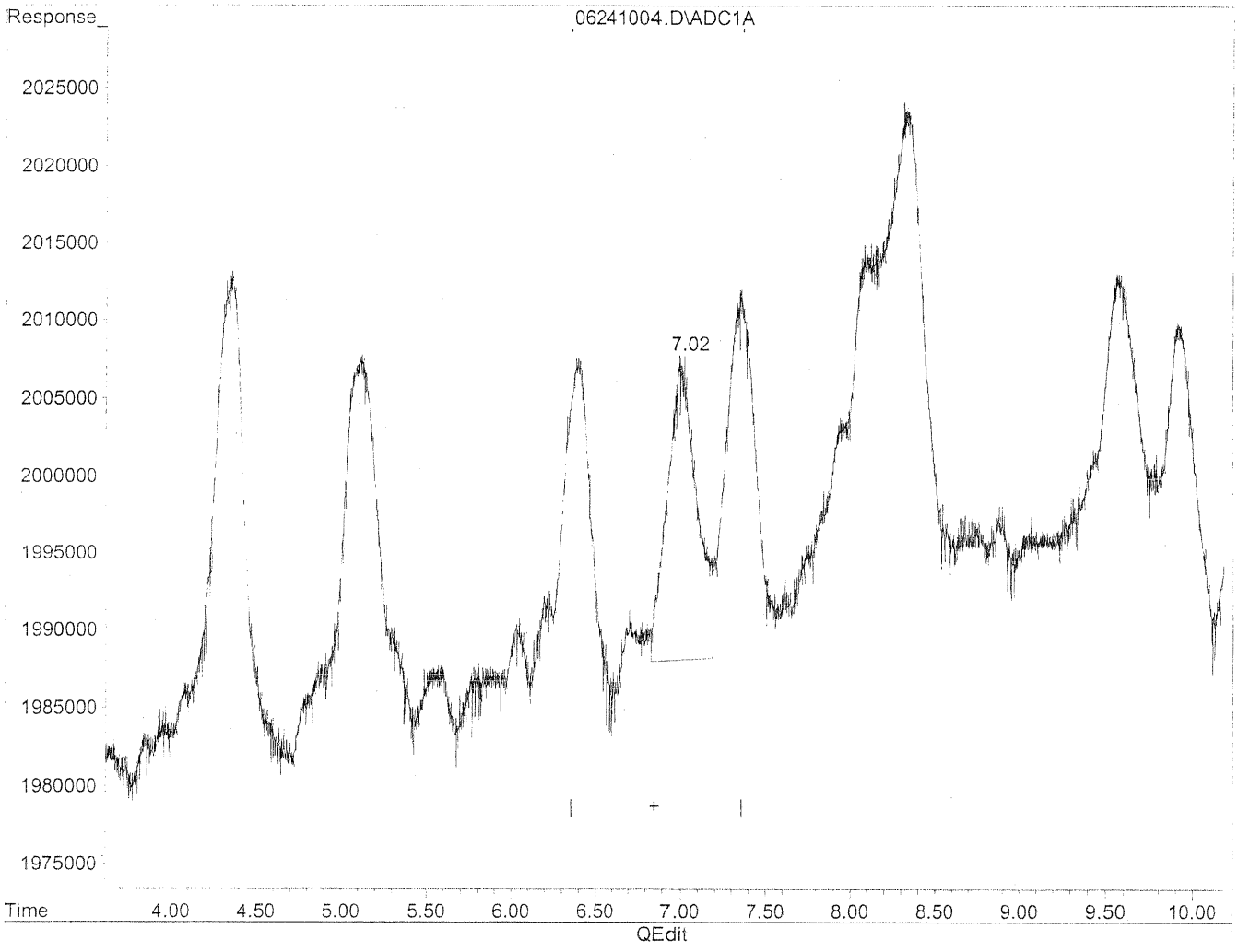
response 2939113

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 16:32:35 2010
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

7.02min 24.701ng/ml m

response 2265328

6/25/10

6/25/10

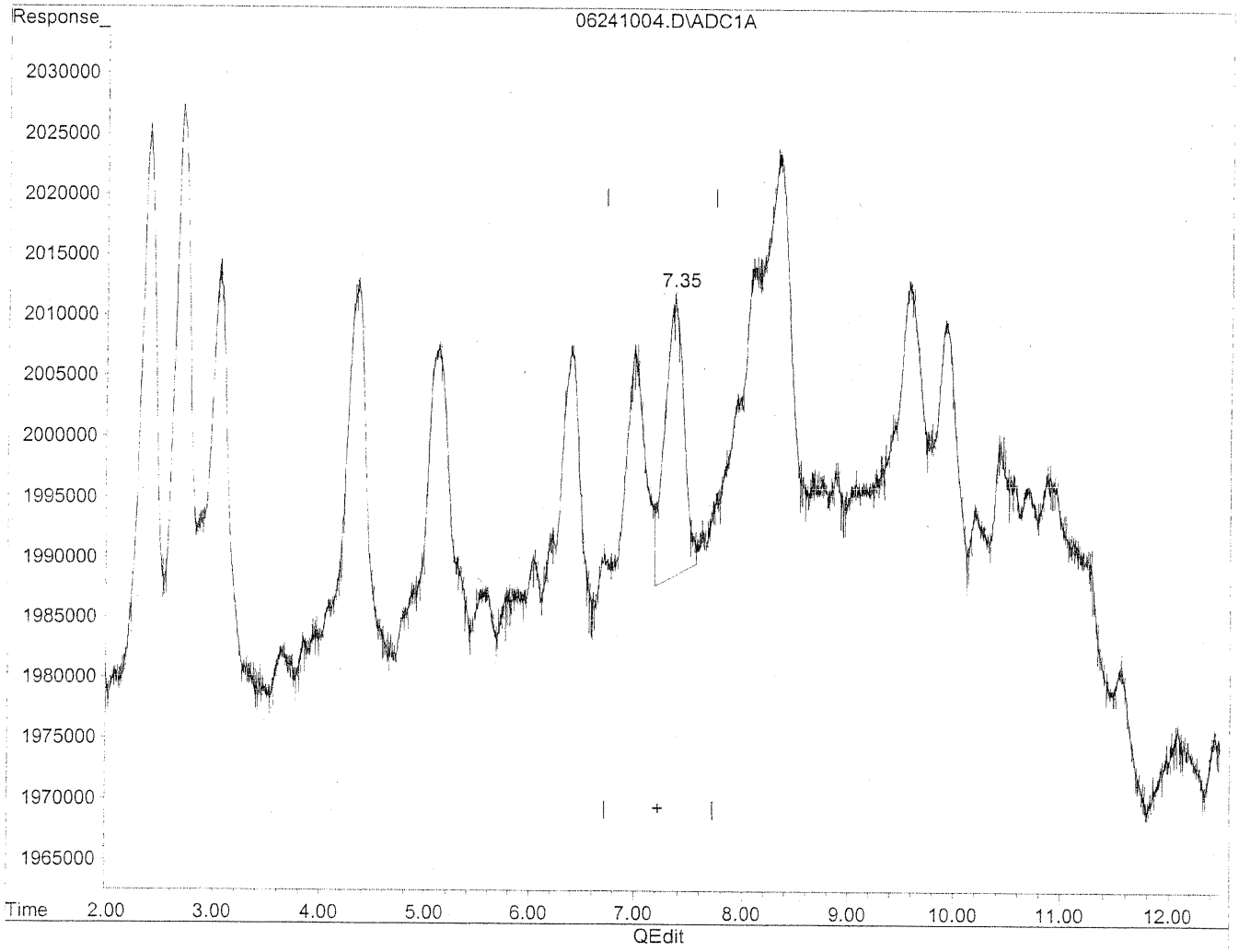
(+) = Expected Retention Time

06241004.D TO110610.M Thu Jun 24 16:34:38 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(8) Valeraldehyde

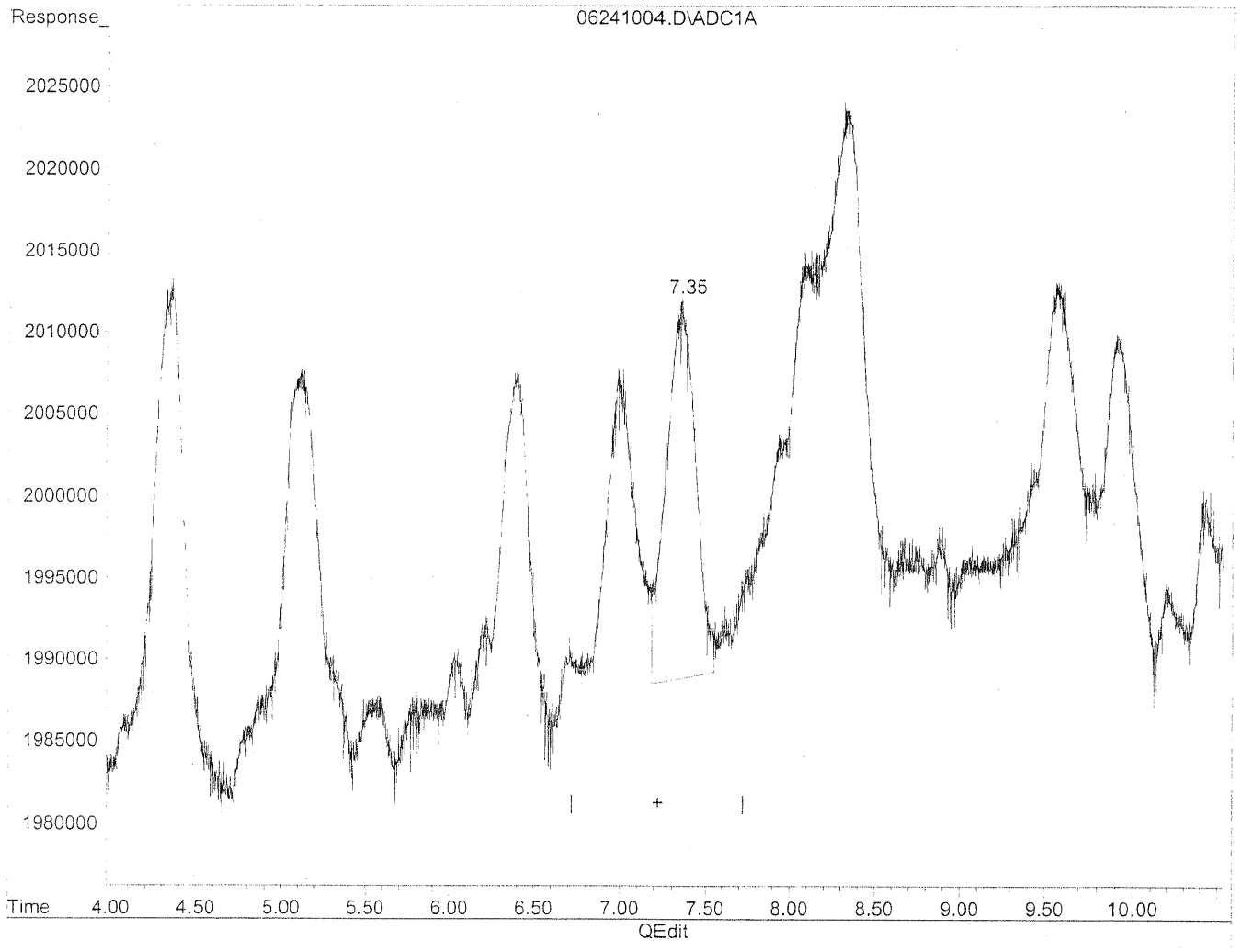
7.35min 32.098ng/ml

response 2777964

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 16:32:35 2010
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.35min 31.385ng/ml m

response 2716276

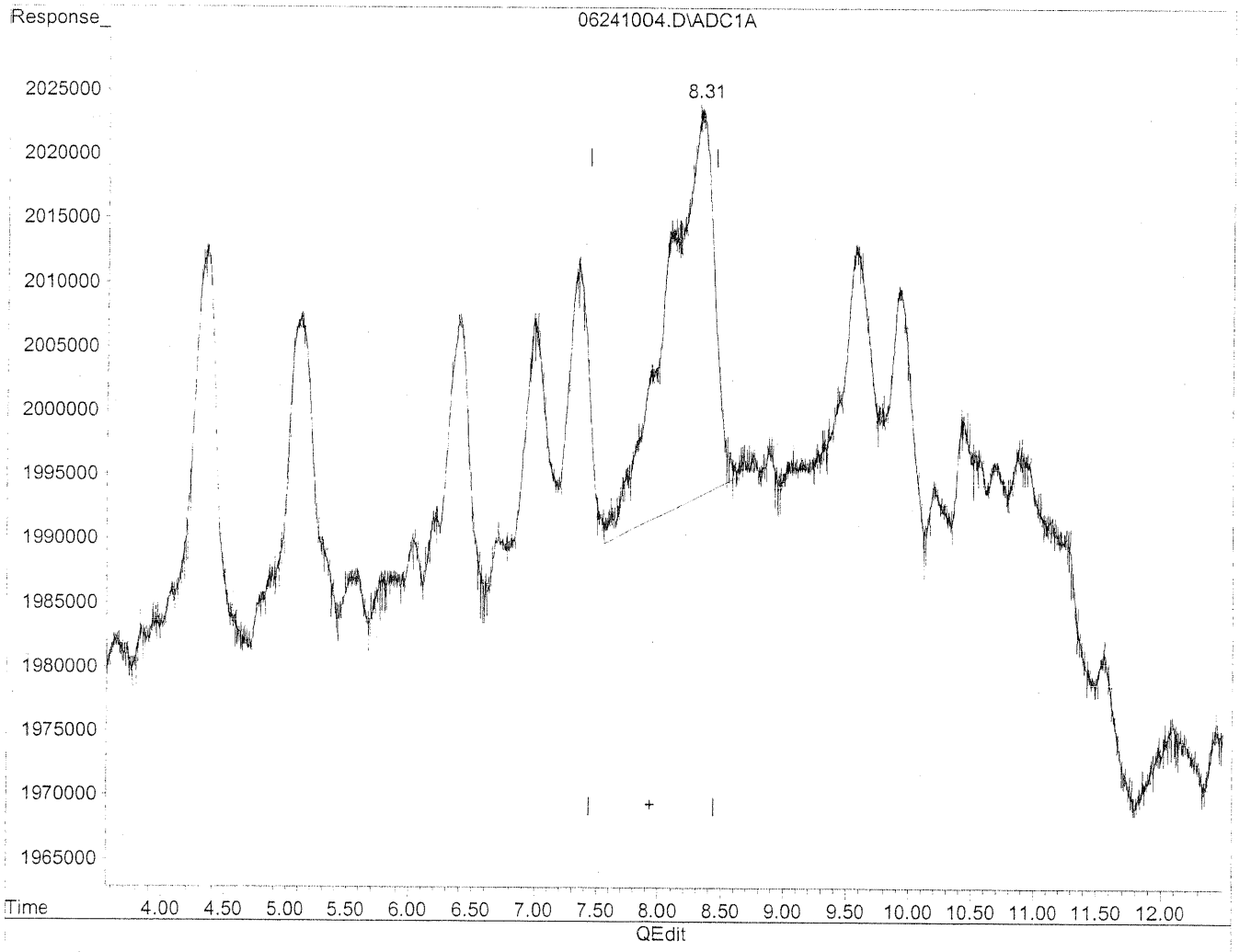
TK 11/28/12

BL 12/25/12

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

8.33min 110.975ng/ml

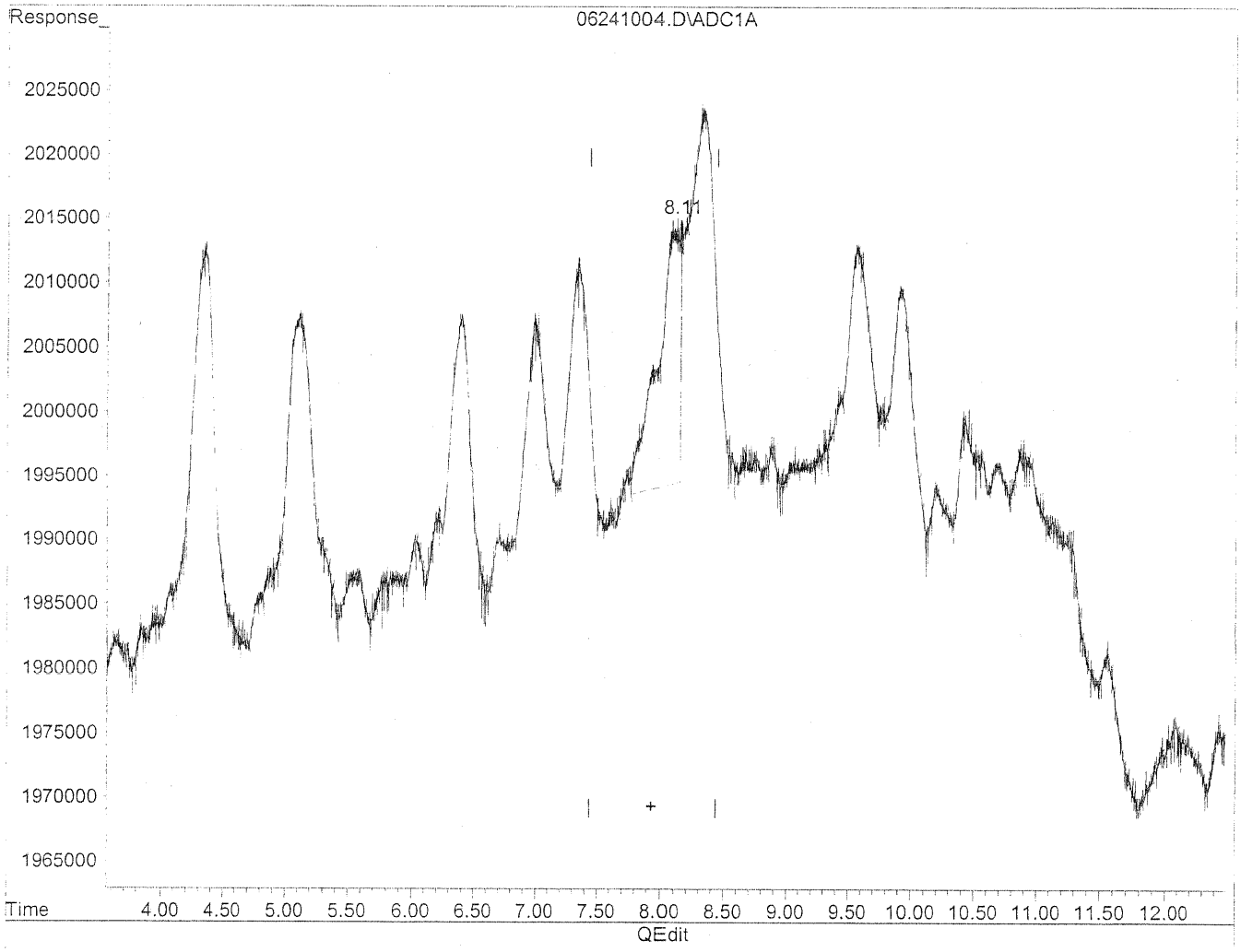
response 7910040

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

8.11min 33.301ng/ml m

response 2373630

6/24/10

6/25/10

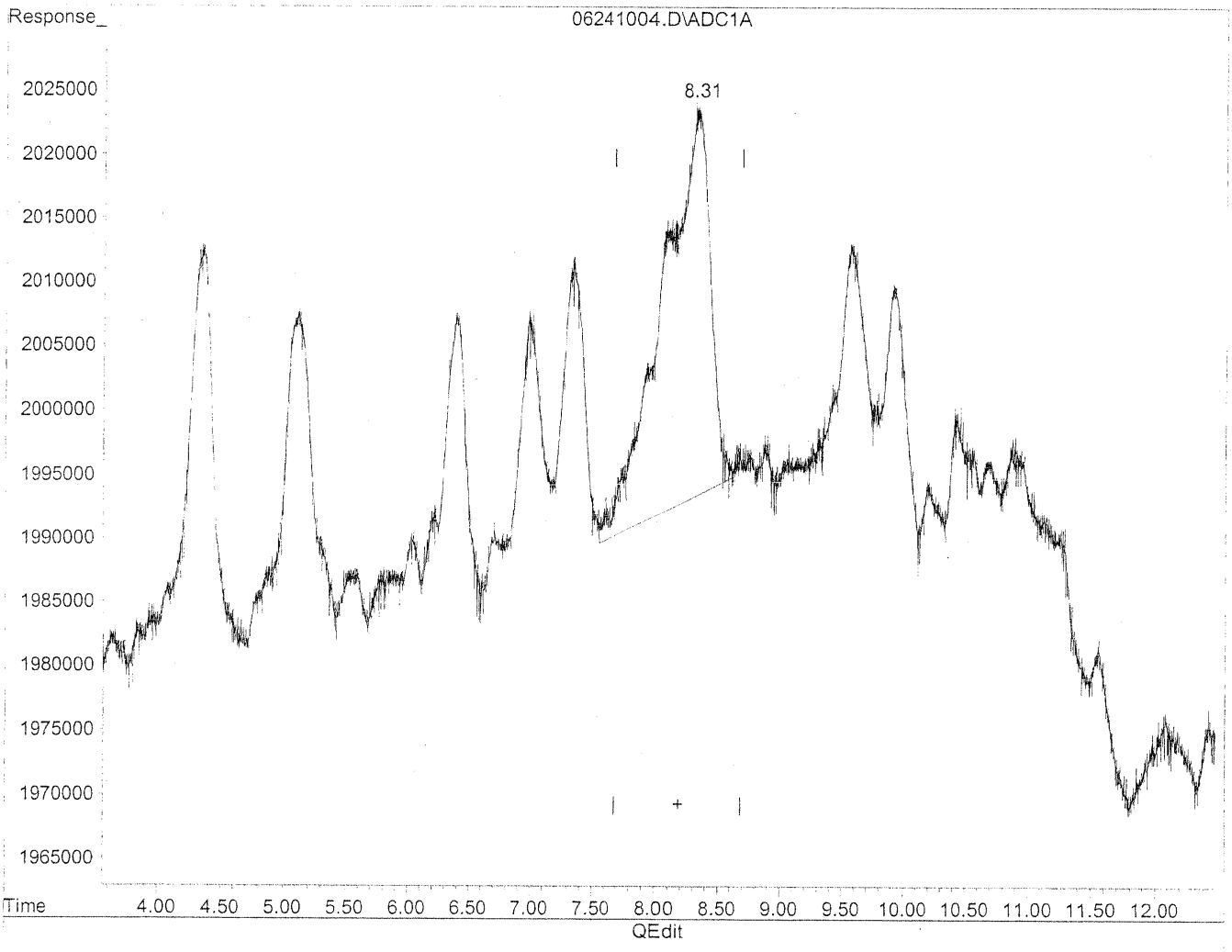
(+) = Expected Retention Time

06241004.D TO110610.M Thu Jun 24 13:29:41 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

8.33min 123.512ng/ml

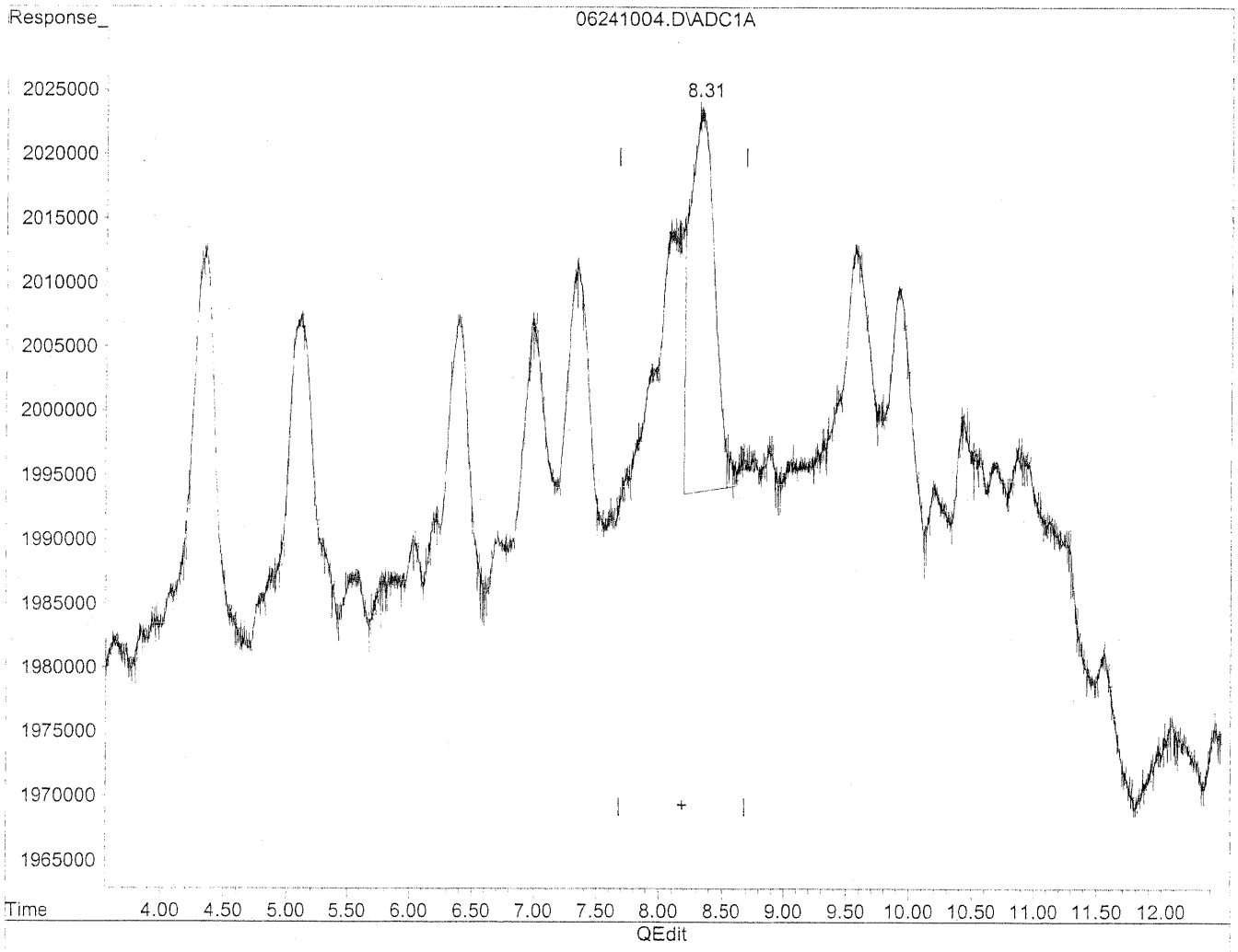
response 7910040

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

8.31min 64.001ng/ml m

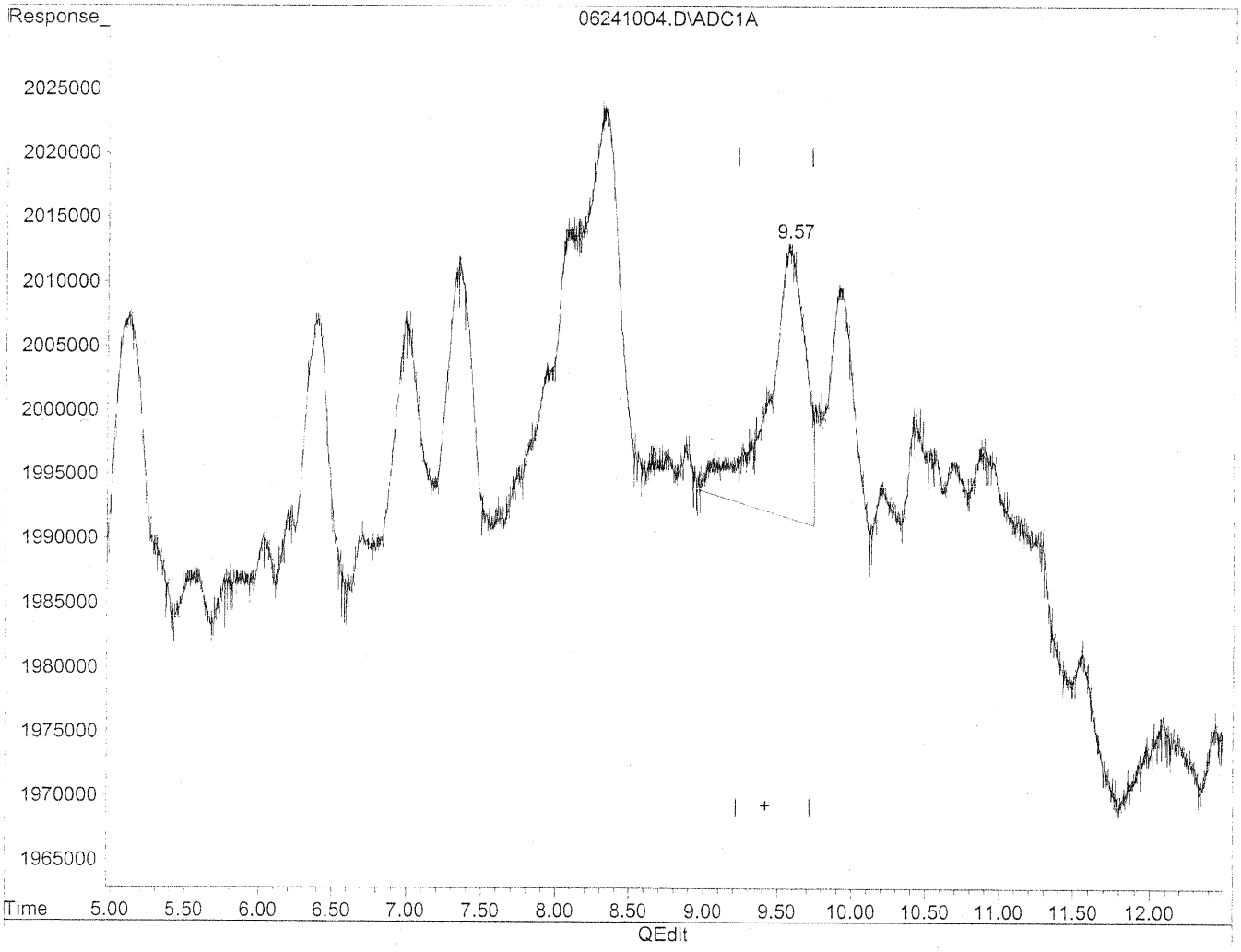
response 4098785

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(11) Hexaldehyde

9.56min 44.941ng/ml

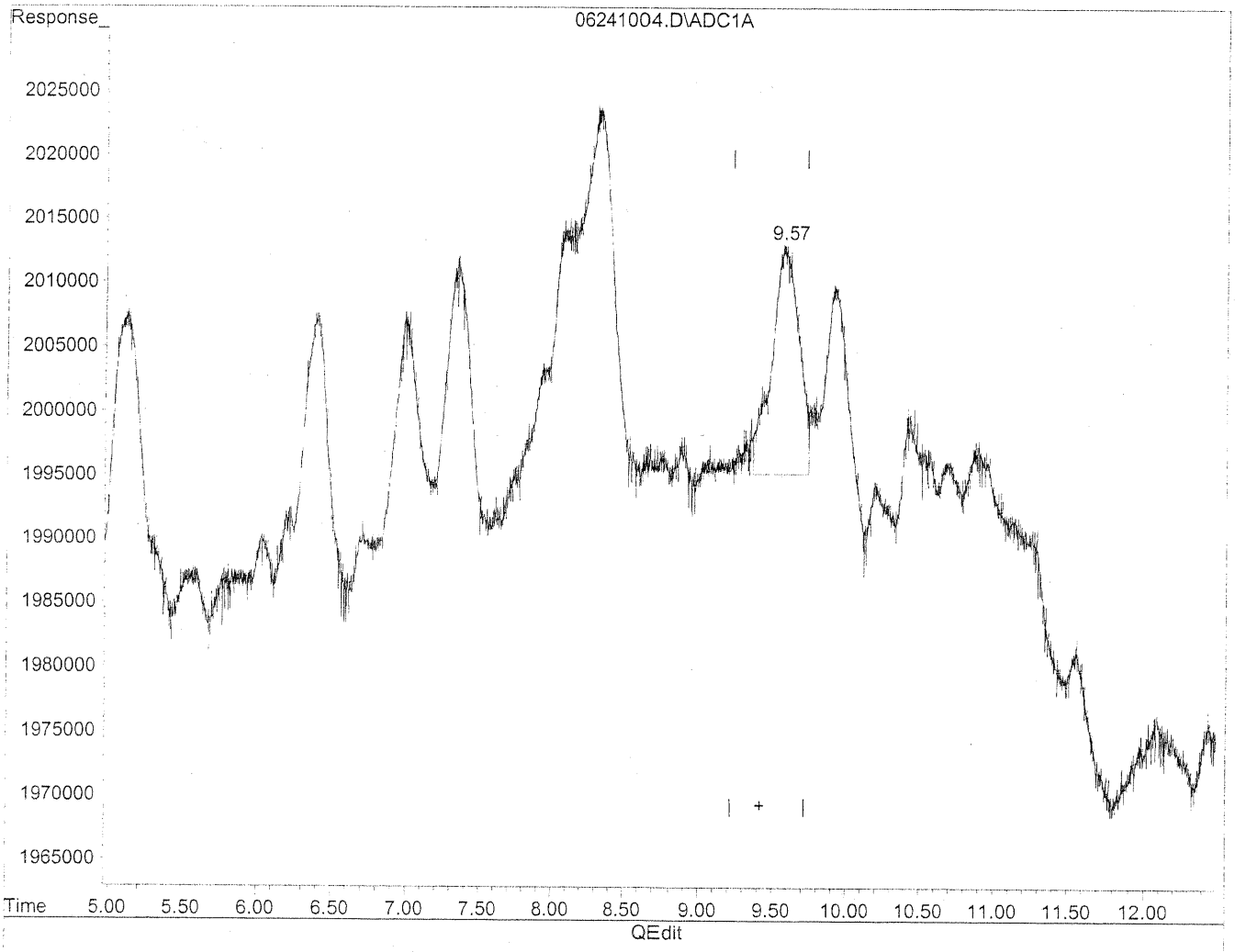
response 3769106

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(11) Hexaldehyde

9.57min 27.309ng/ml m

response 2290365

HC 6/28/10

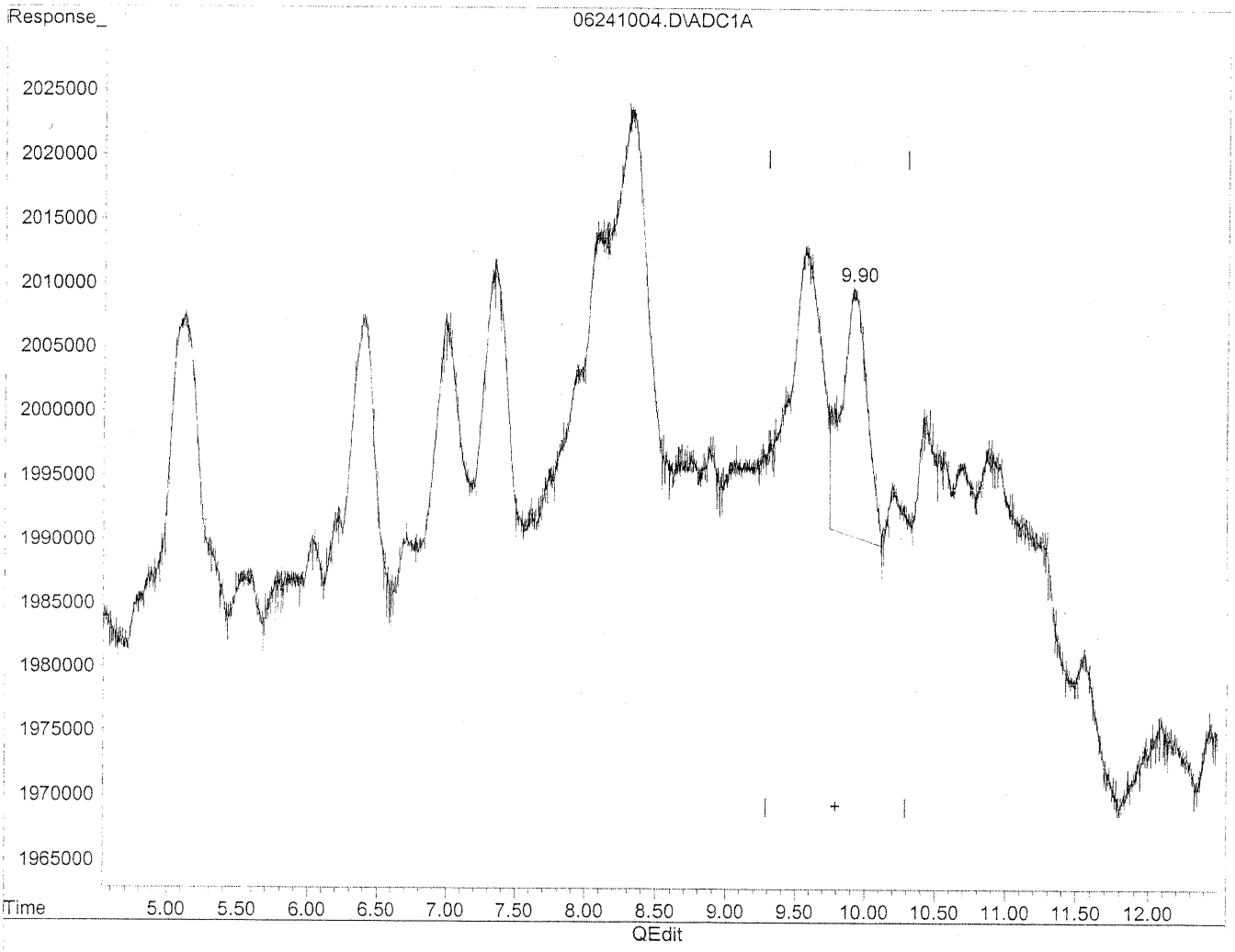
6/25/10

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:26 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

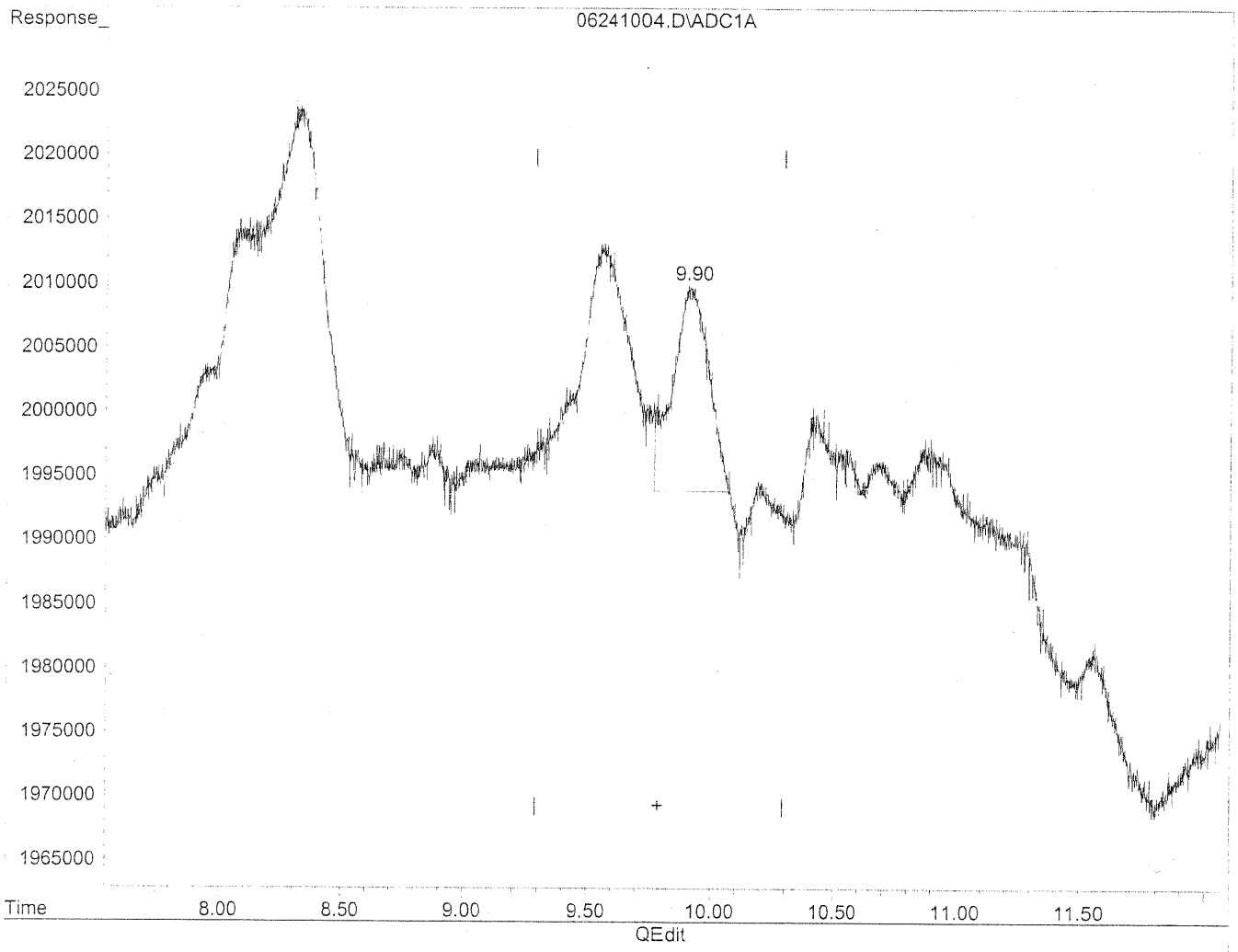
9.91min 43.373ng/ml

response 2473055

(+) = Expected Retention Time

Data File : J:\LC01\DATA\TO11\2010_06\24\06241004.D Vial: 2
Acq On : 24 Jun 2010 12:04 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:41 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 16:42:47 2010
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

9.90min 28.403ng/ml m

response 1619465

TC 6/25/10

TC 6/25/10

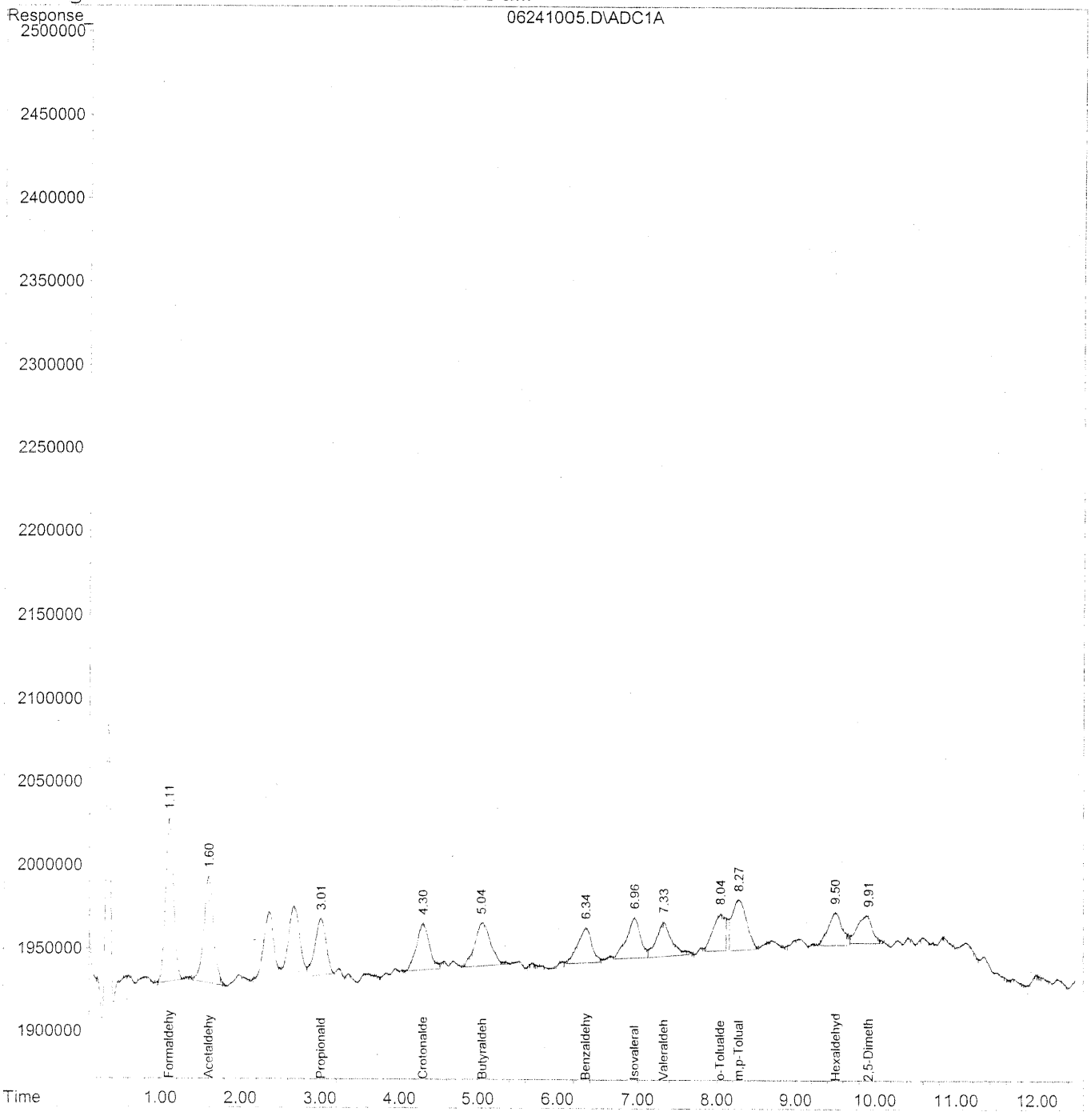
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:43 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
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\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:43 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Initial Calibration
DataAcq Meth : TO11S.M

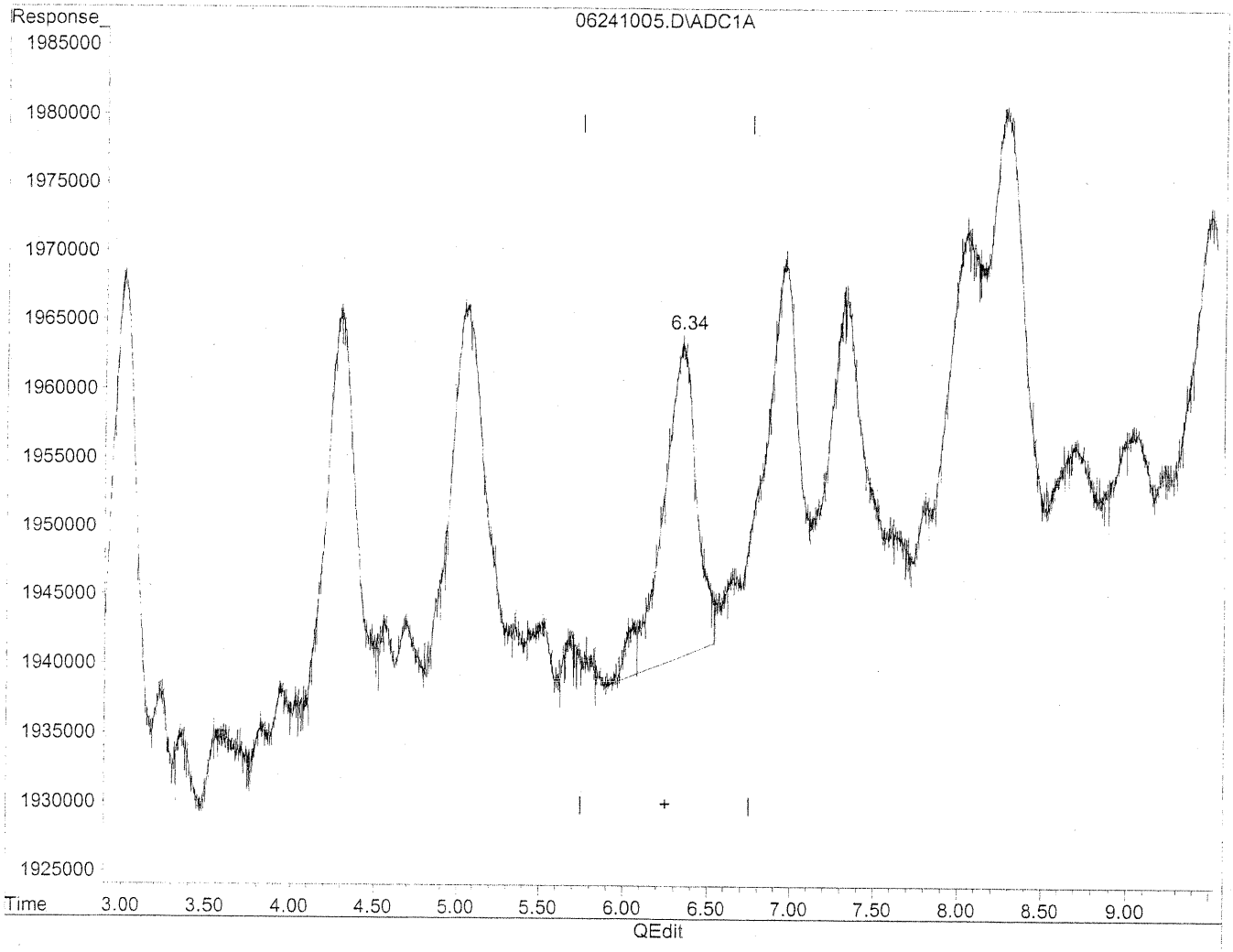
Volume Inj. : 5uL
Signal Phase : Supleco Supelcosil LC-18
Signal Info : 3.3cm x 3mm x 3um

	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.11	6598476	32.263	ng/ml
2)	Acetaldehyde	1.60	5336318	34.153	ng/ml
3)	Propionaldehyde	3.01	3364453	28.440	ng/ml
4)	Crotonaldehyde	4.30	3240110	30.081	ng/ml
5)	Butyraldehyde	5.05	3652578	36.738	ng/ml
6)	Benzaldehyde	6.34	2527622	35.282	ng/mlm
7)	Isovaleraldehyde	6.96	2995744	36.116	ng/mlm
8)	Valeraldehyde	7.33	2768346	34.756	ng/mlm
9)	o-Tolualdehyde	8.04	2334081	35.419	ng/mlm
10)	m,p-Tolualdehyde	8.27	3866830	64.315	ng/mlm
11)	Hexaldehyde	9.51	2424895	32.009	ng/mlm
12)	2,5-Dimethylbenzaldehyde	9.91	2173046	41.846	ng/mlm

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.34min 44.427ng/ml

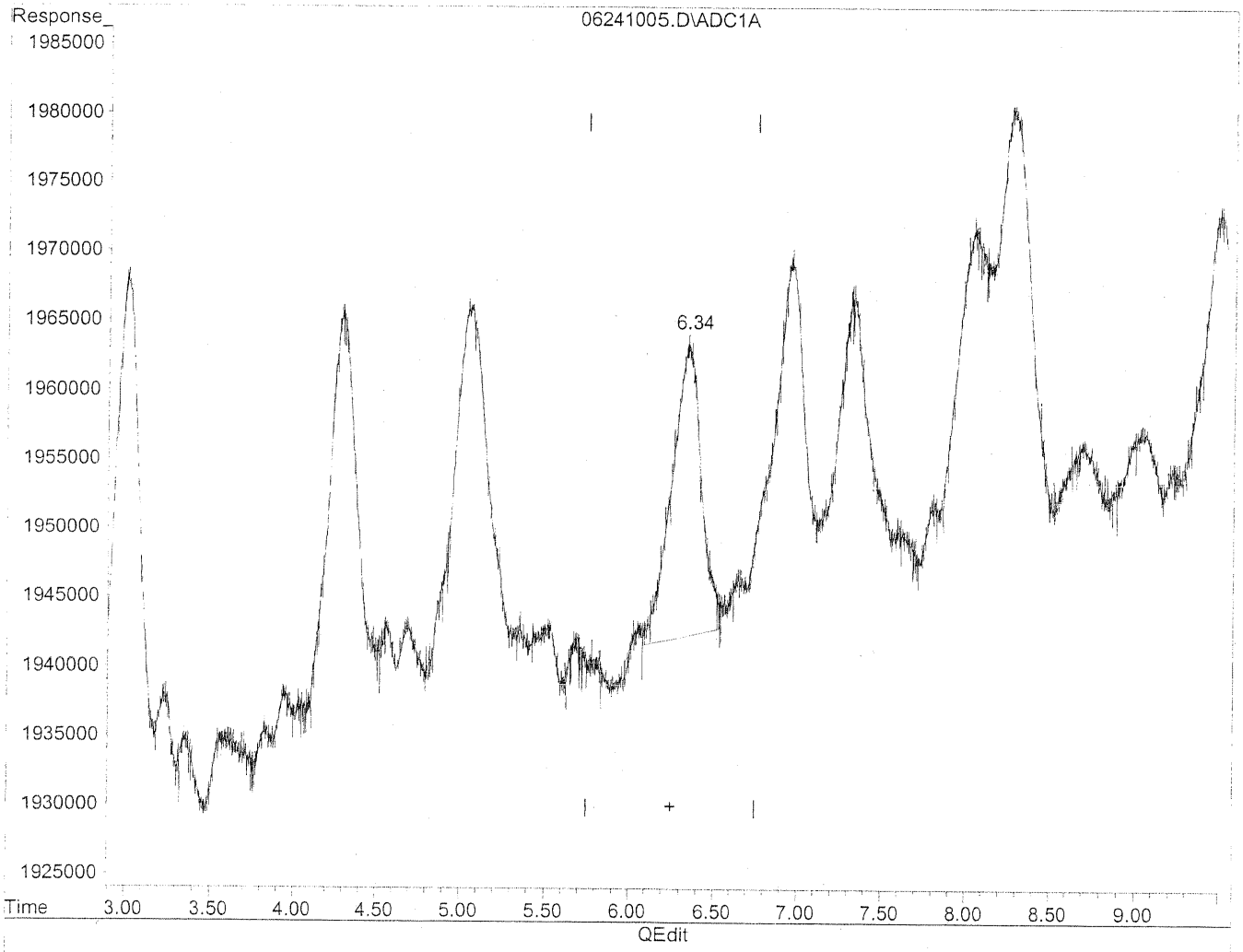
response 3182803

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.34min 35.282ng/ml m

response 2527622

HC 6/28/10

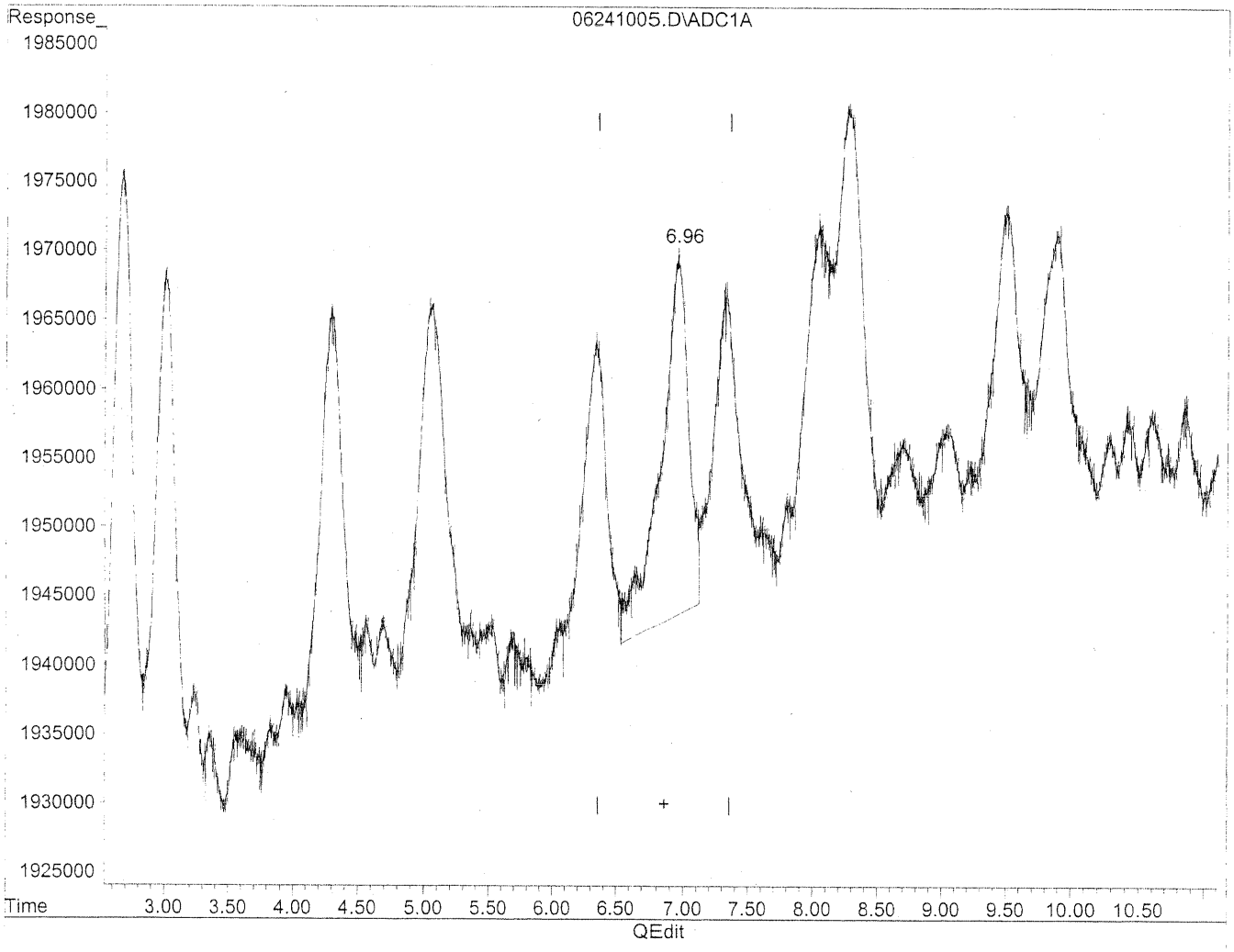
TC 6/25/10

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

6.96min 45.094ng/ml

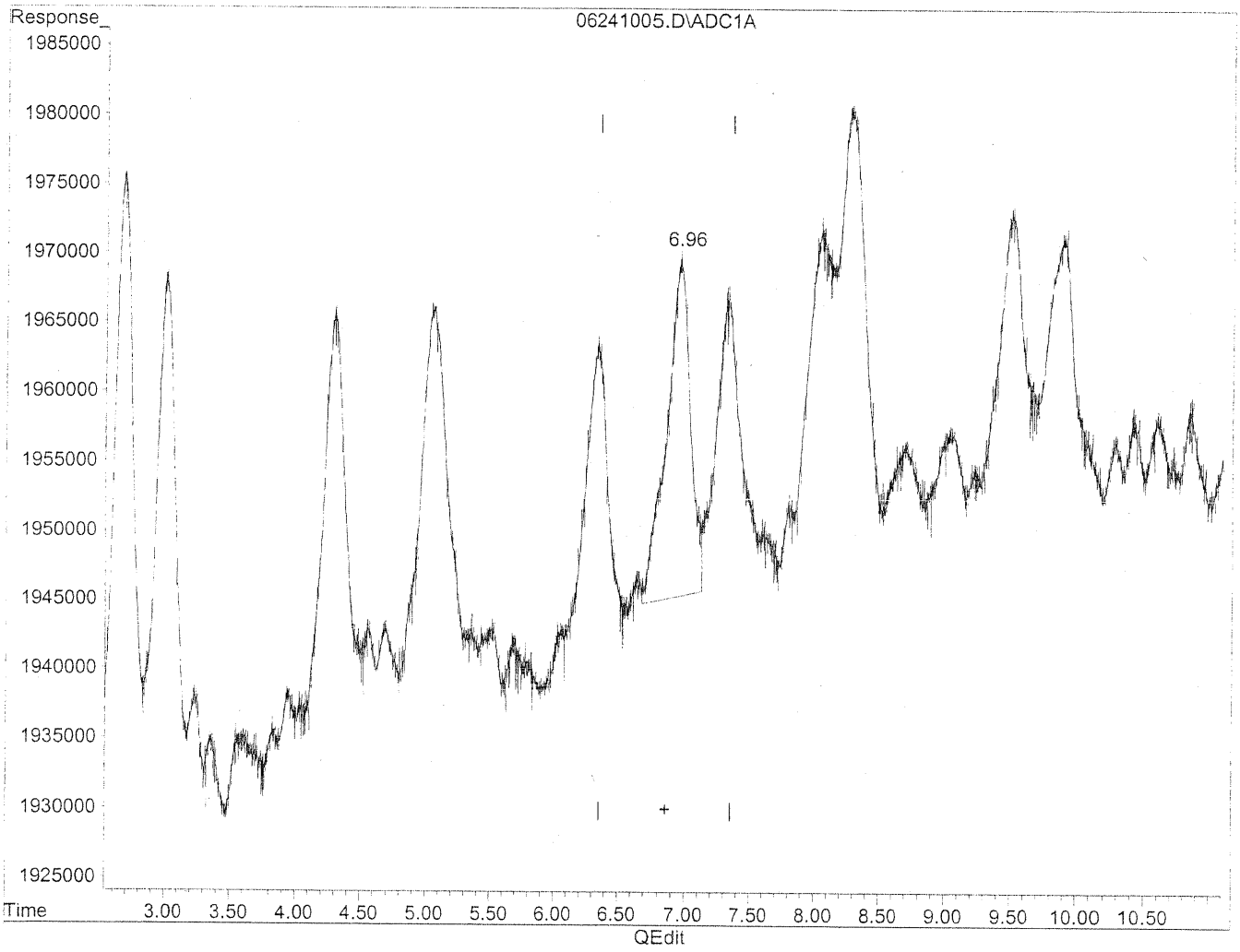
response 3740491

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

6.96min 36.116ng/ml m

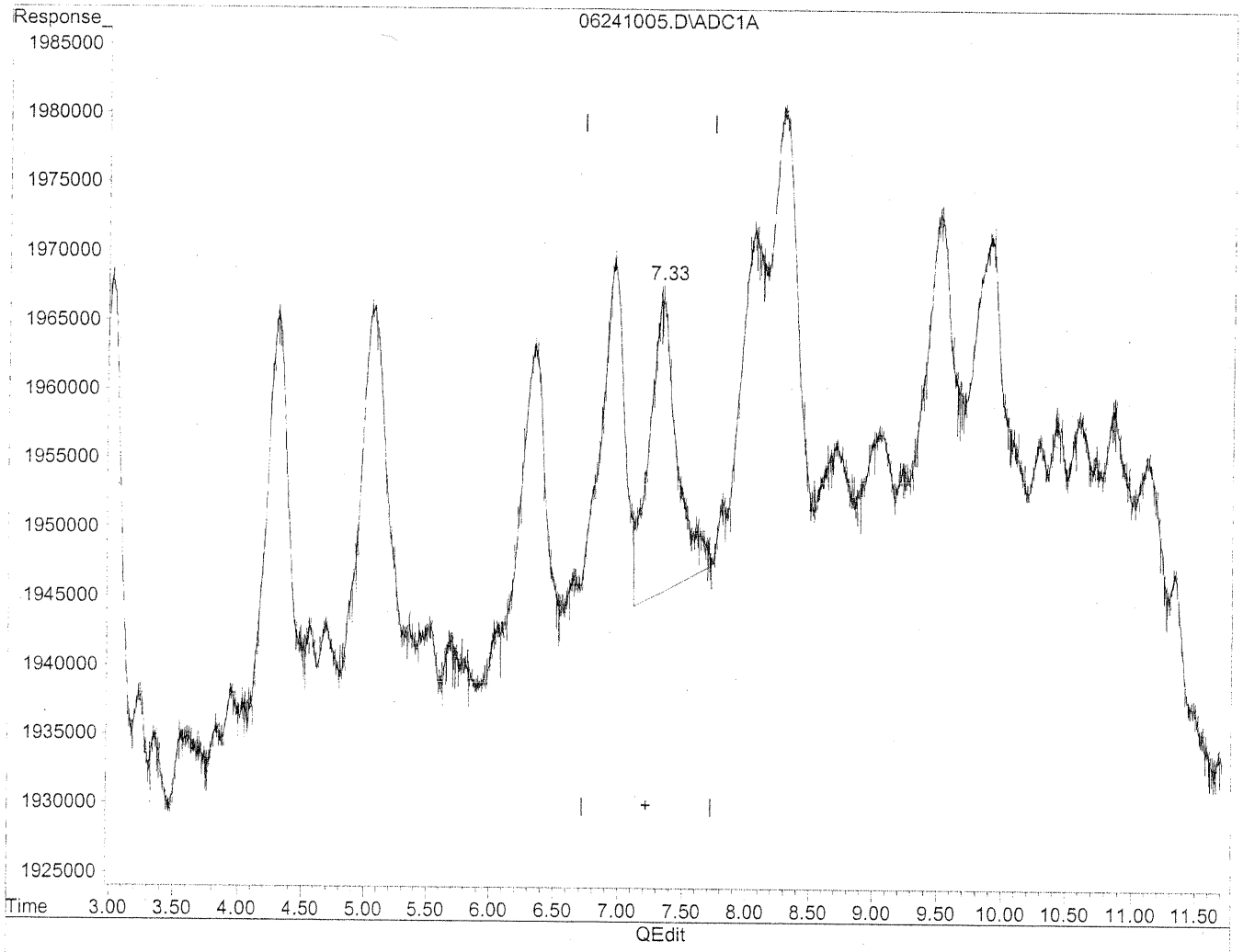
response 2995744

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.33min 39.386ng/ml

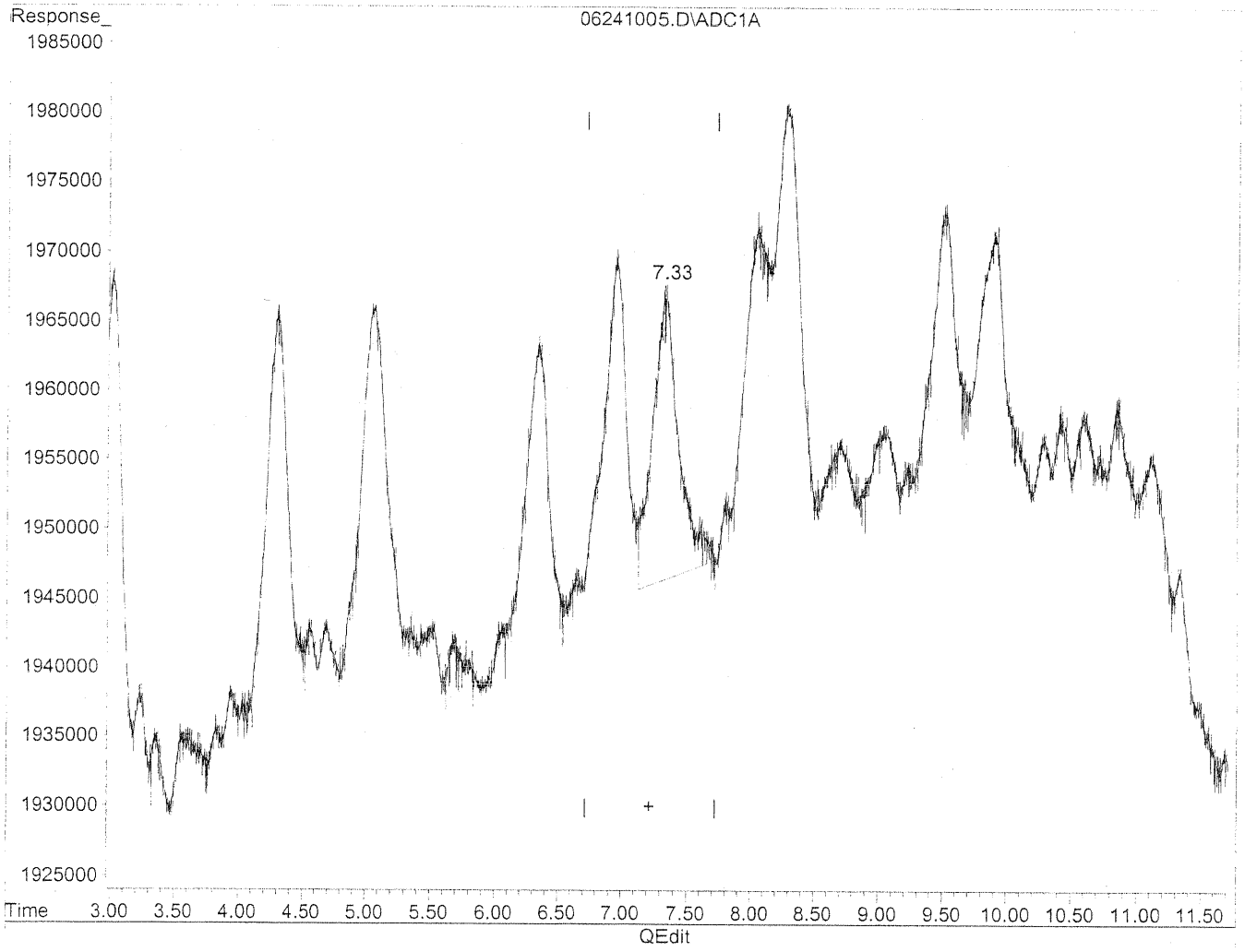
response 3137126

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.33min 34.756ng/ml m

response 2768346

HC 1/28/10

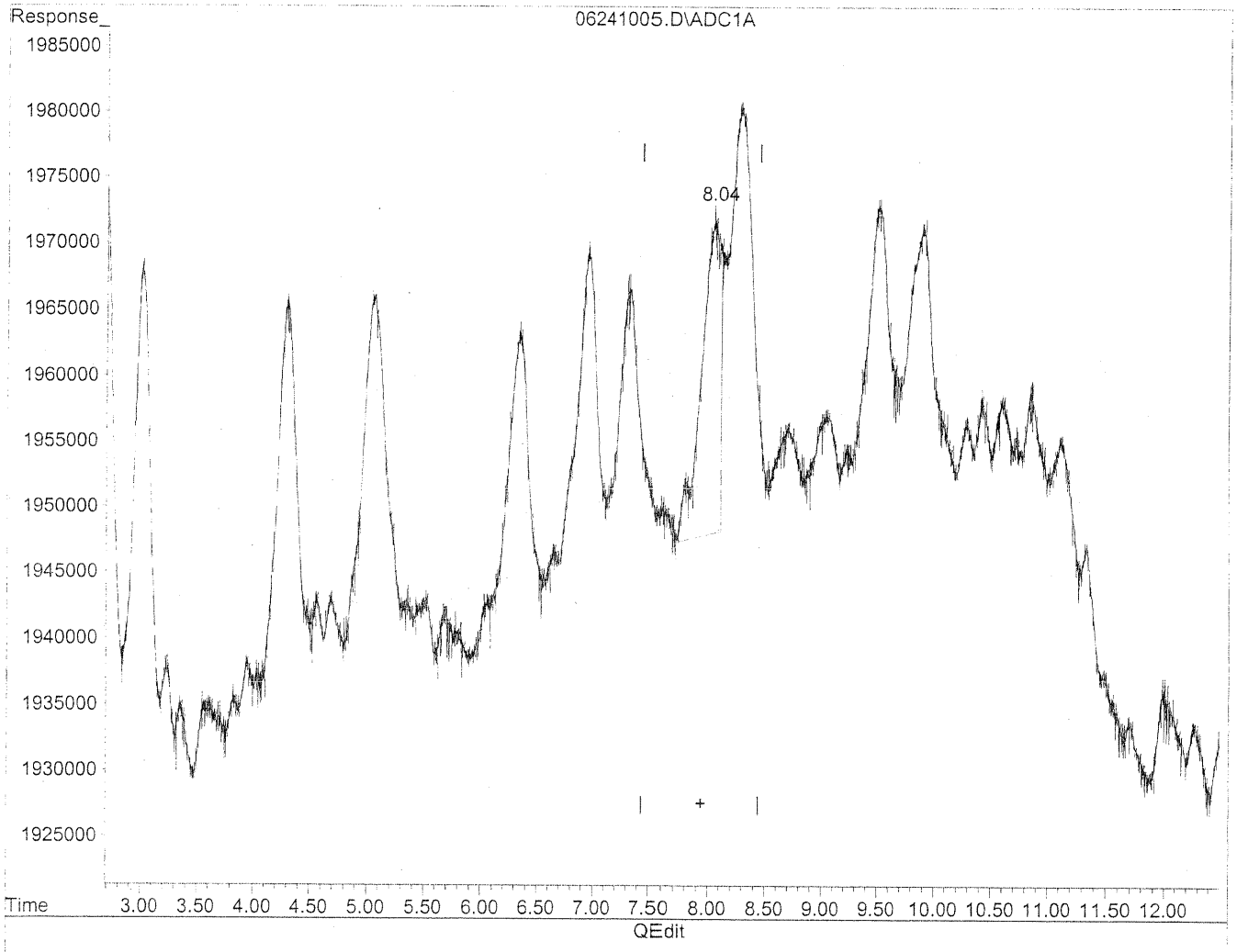
TC 6/28/10

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

8.04min 41.808ng/ml

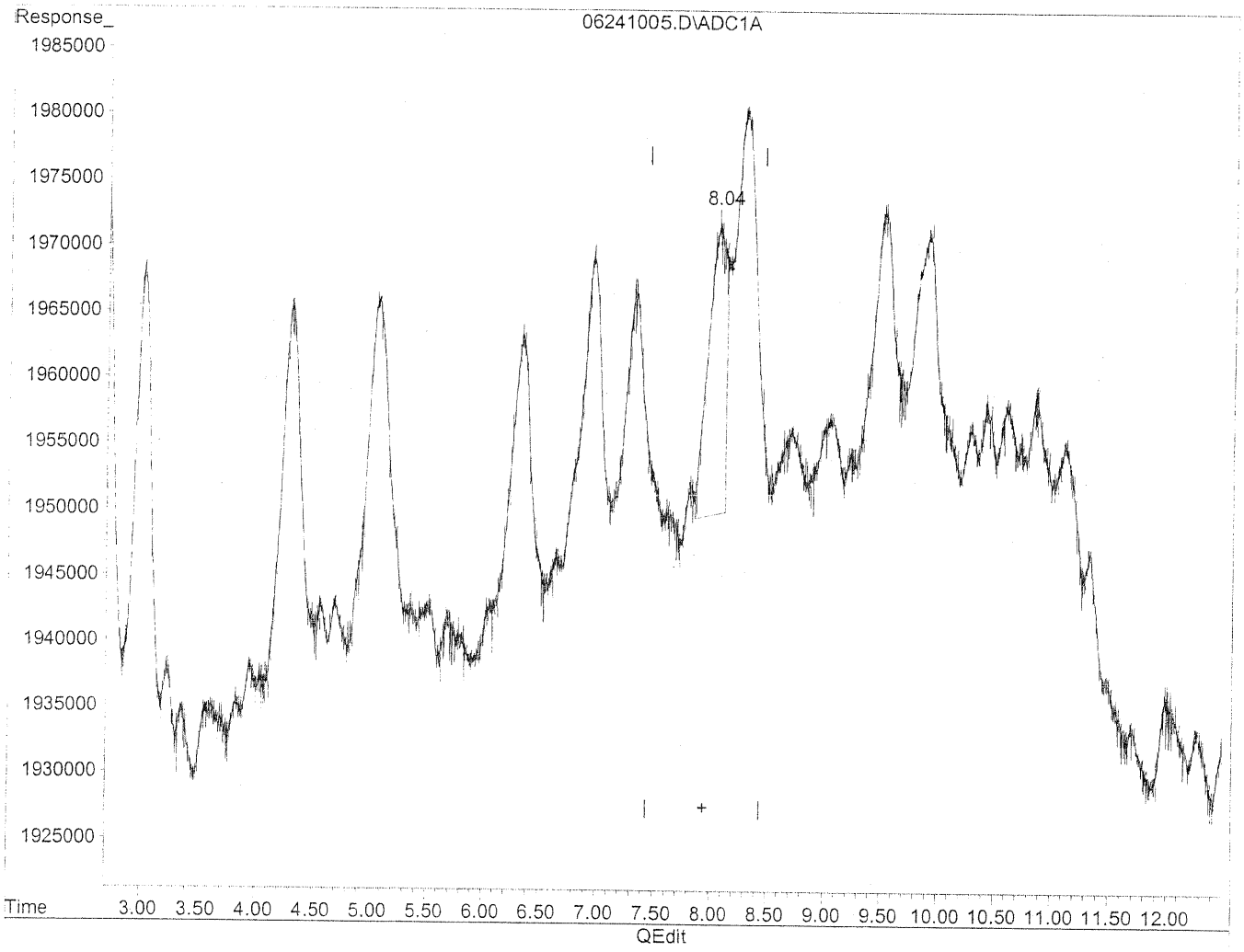
response 2755141

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

8.04min 35.419ng/ml m

response 2334081

HC 6/25/10

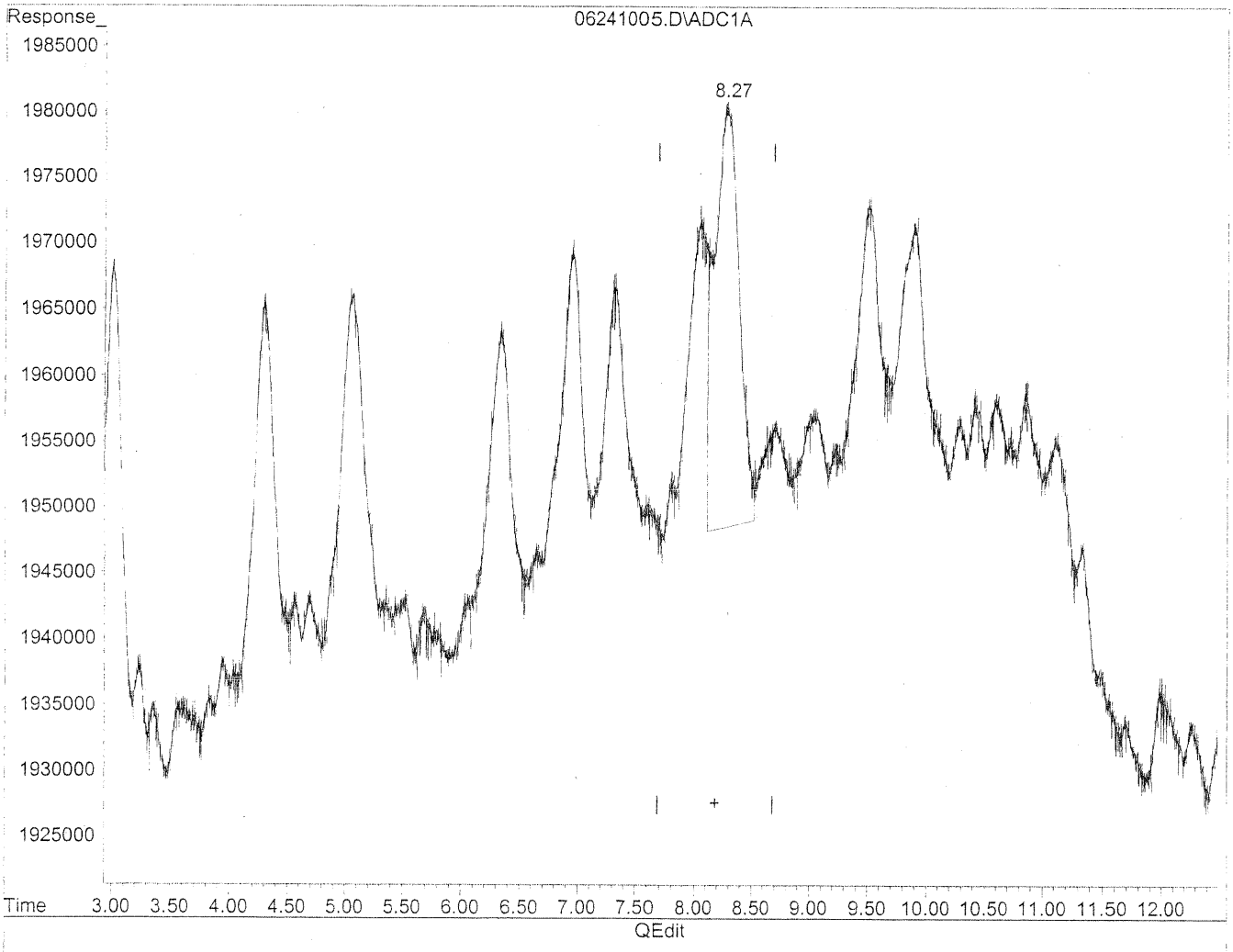
Sh, BC 6/28/10

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

8.27min 79.685ng/ml

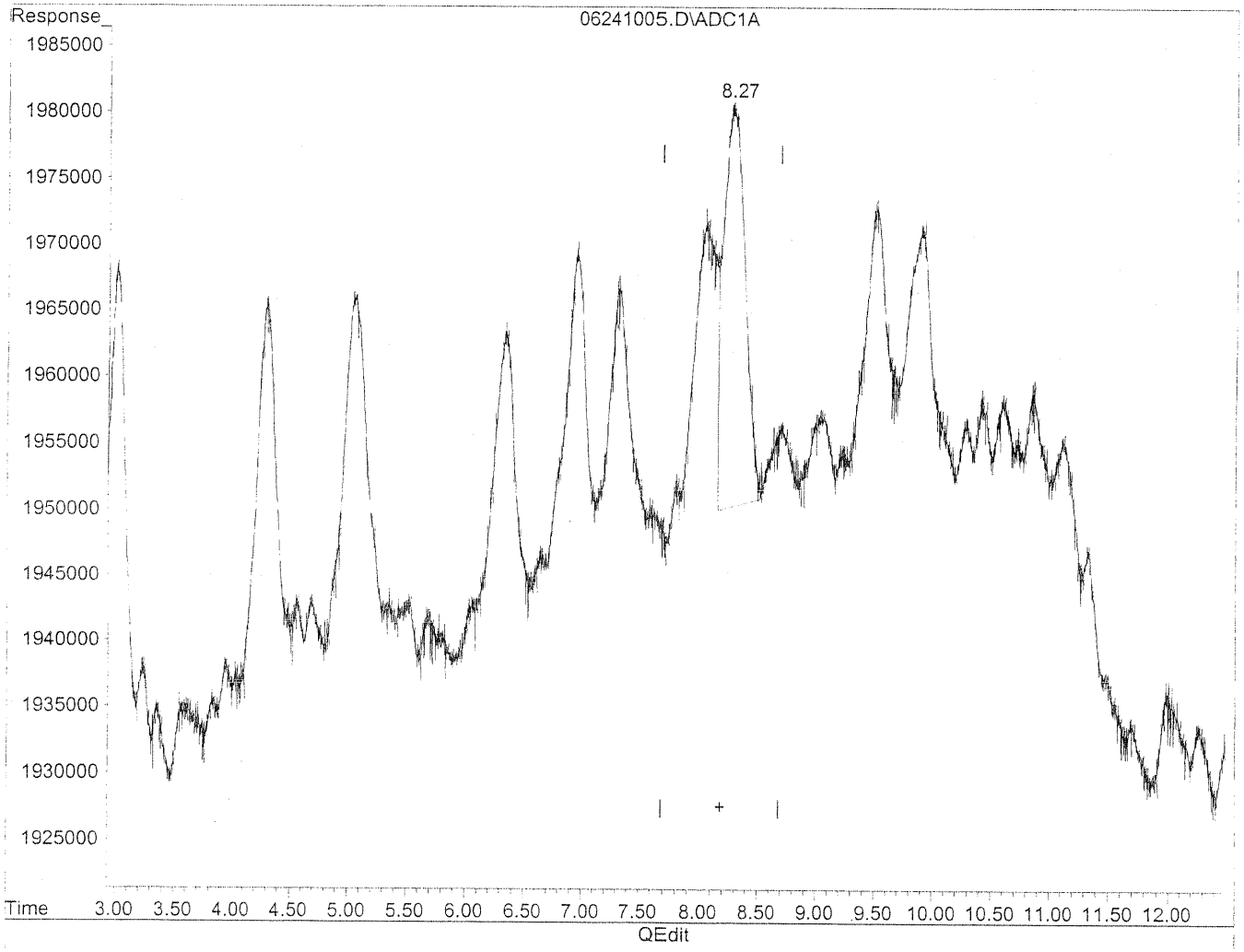
response 4790889

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

8.27min 64.315ng/ml m

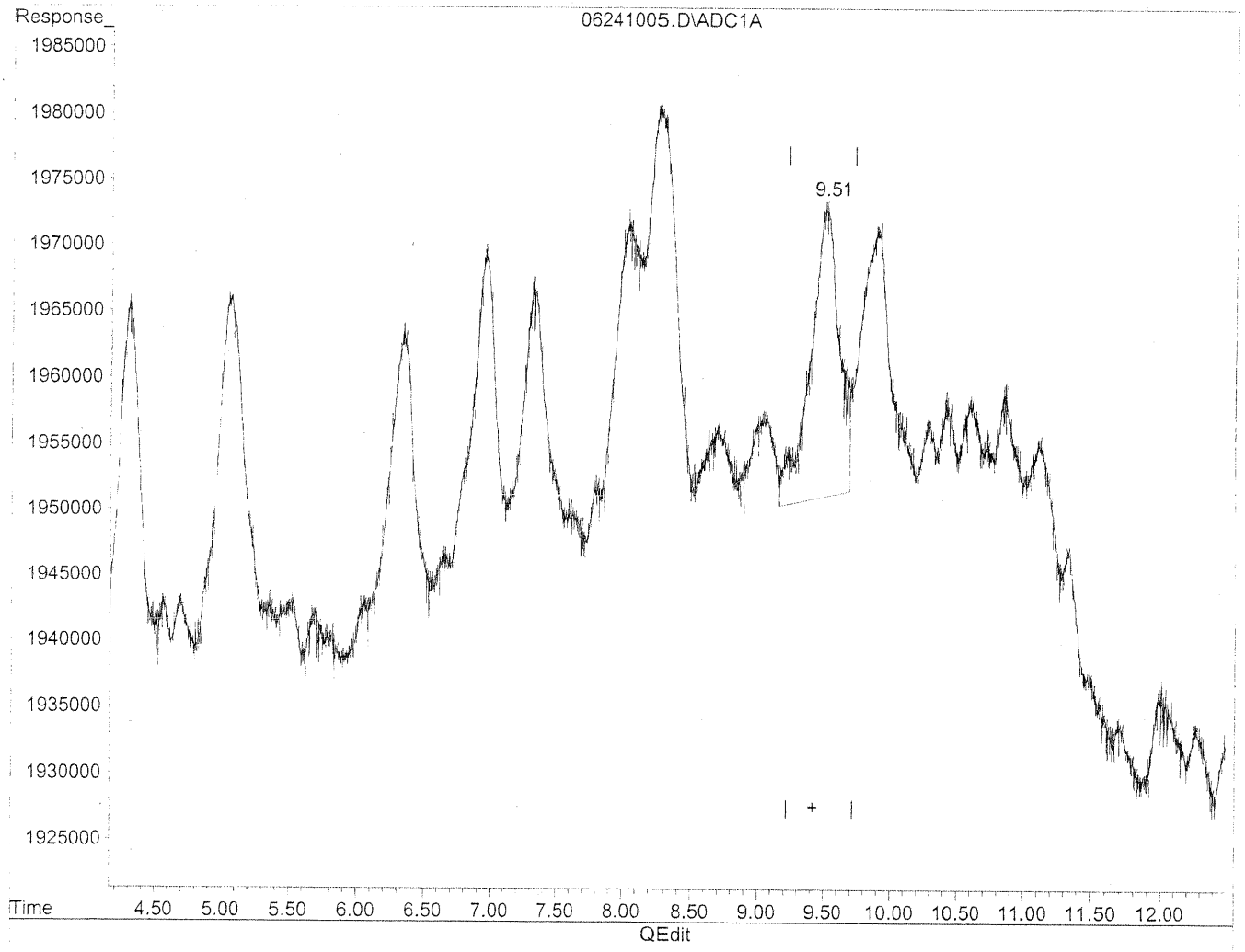
response 3866830

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(11) Hexaldehyde

9.50min 43.500ng/ml

response 3295400

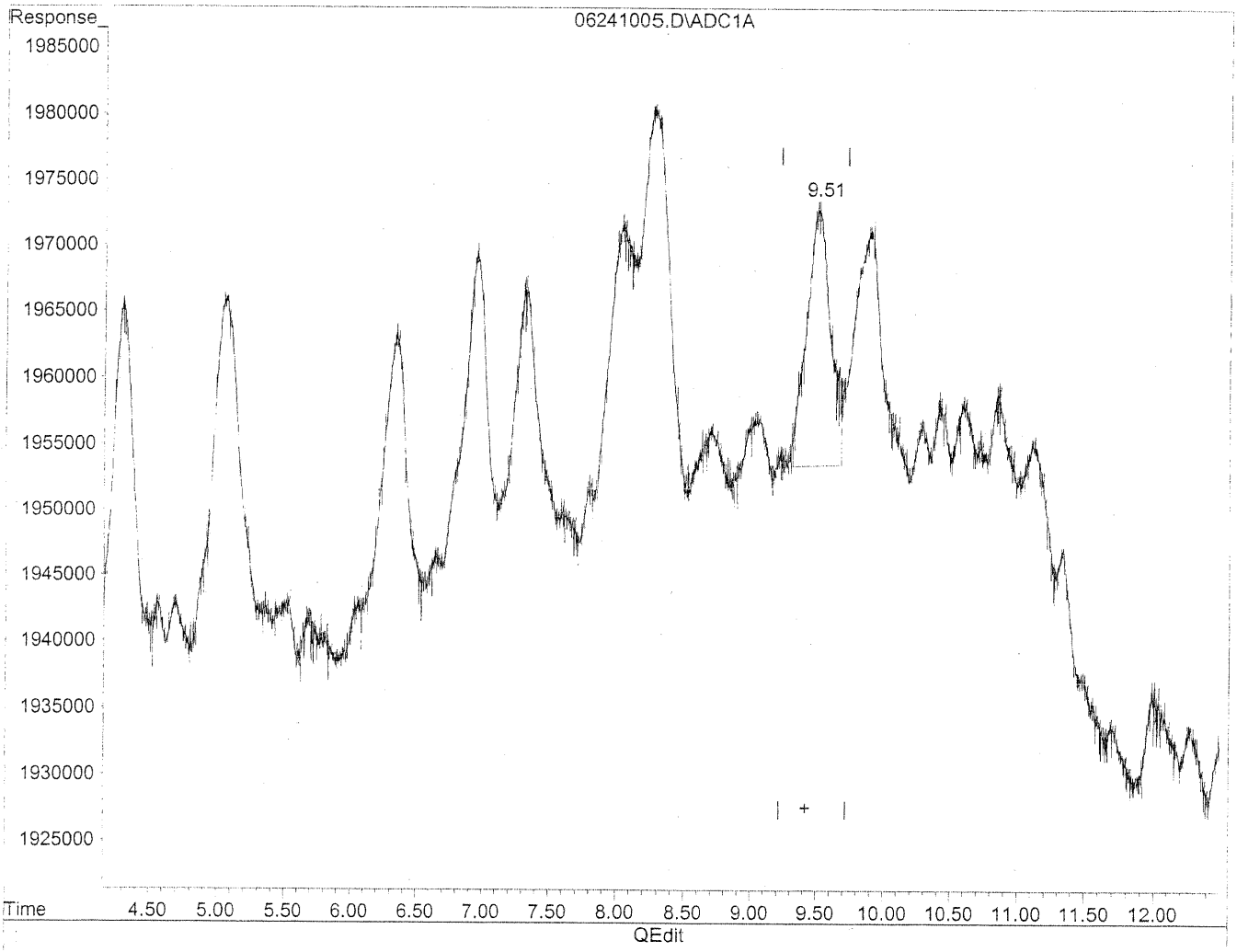
(+) = Expected Retention Time

06241005.D TO110610.M Thu Jun 24 13:32:28 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(11) Hexaldehyde

9.51min 32.009ng/ml m

response 2424895

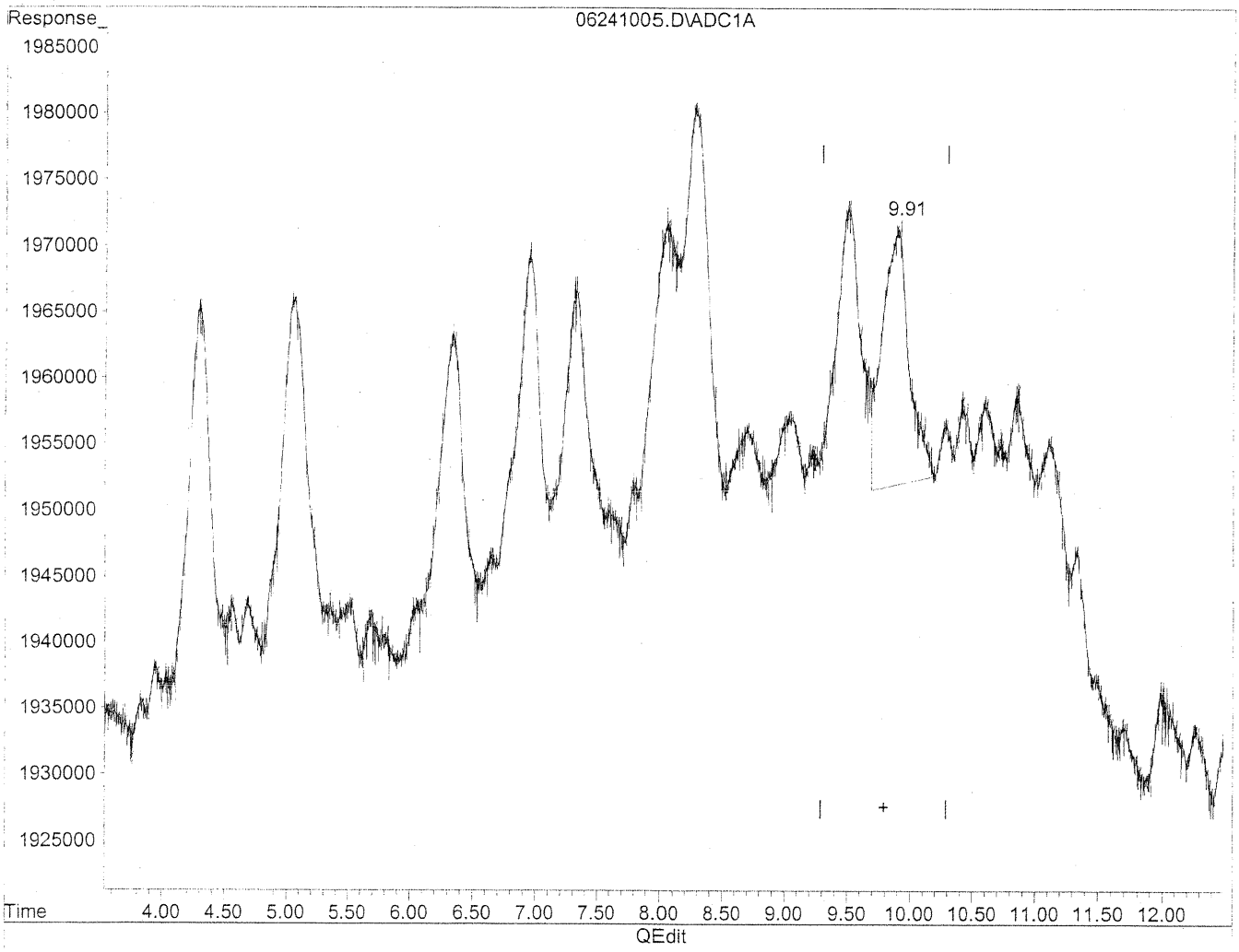
(+) = Expected Retention Time

06241005.D TO110610.M Thu Jun 24 13:32:43 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:30 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

9.89min 56.170ng/ml

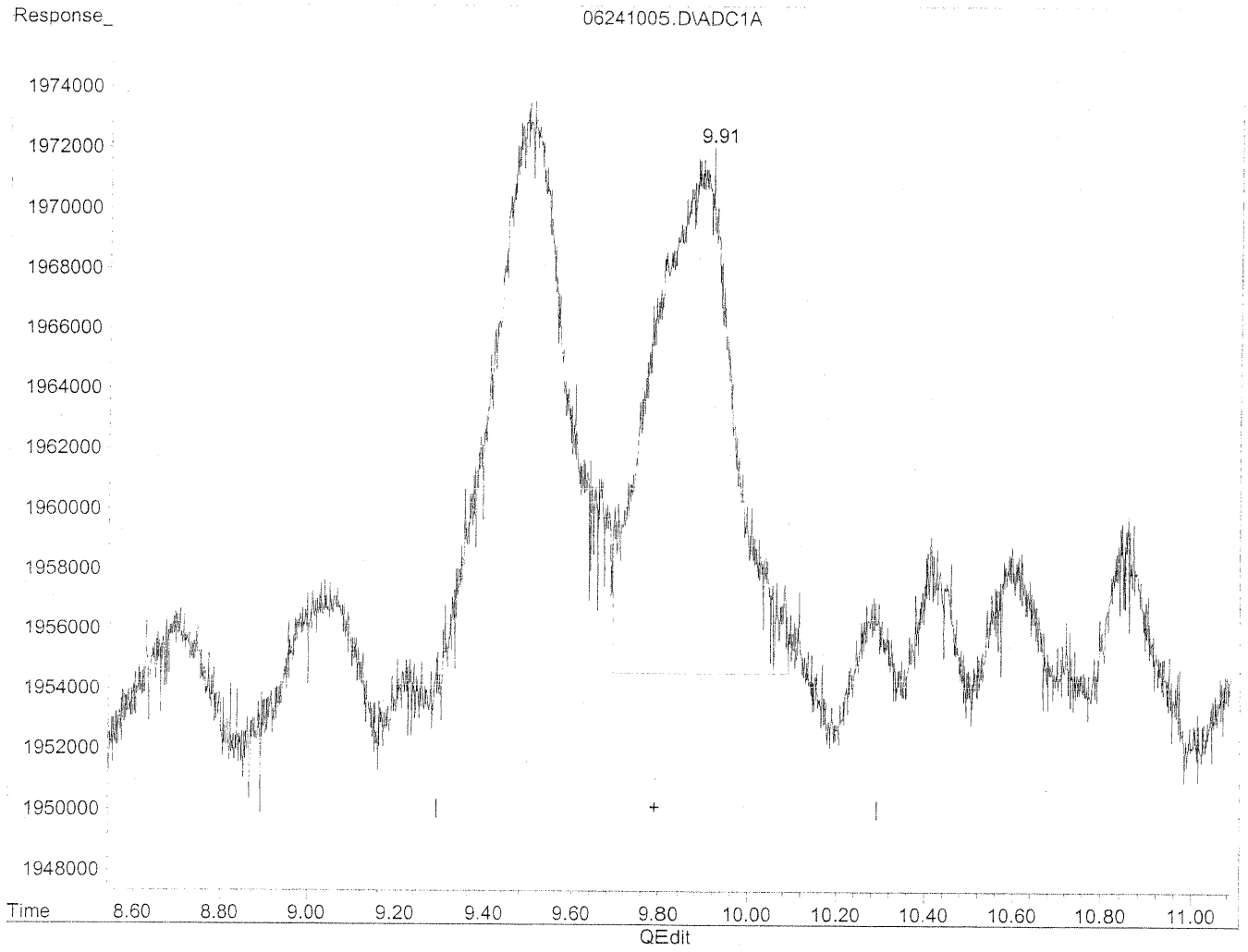
response 2916898

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241005.D Vial: 3
Acq On : 24 Jun 2010 12:18 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:43 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 16:42:47 2010
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

9.91min 41.846ng/ml m

response 2173046

6/25/10

6/25/10

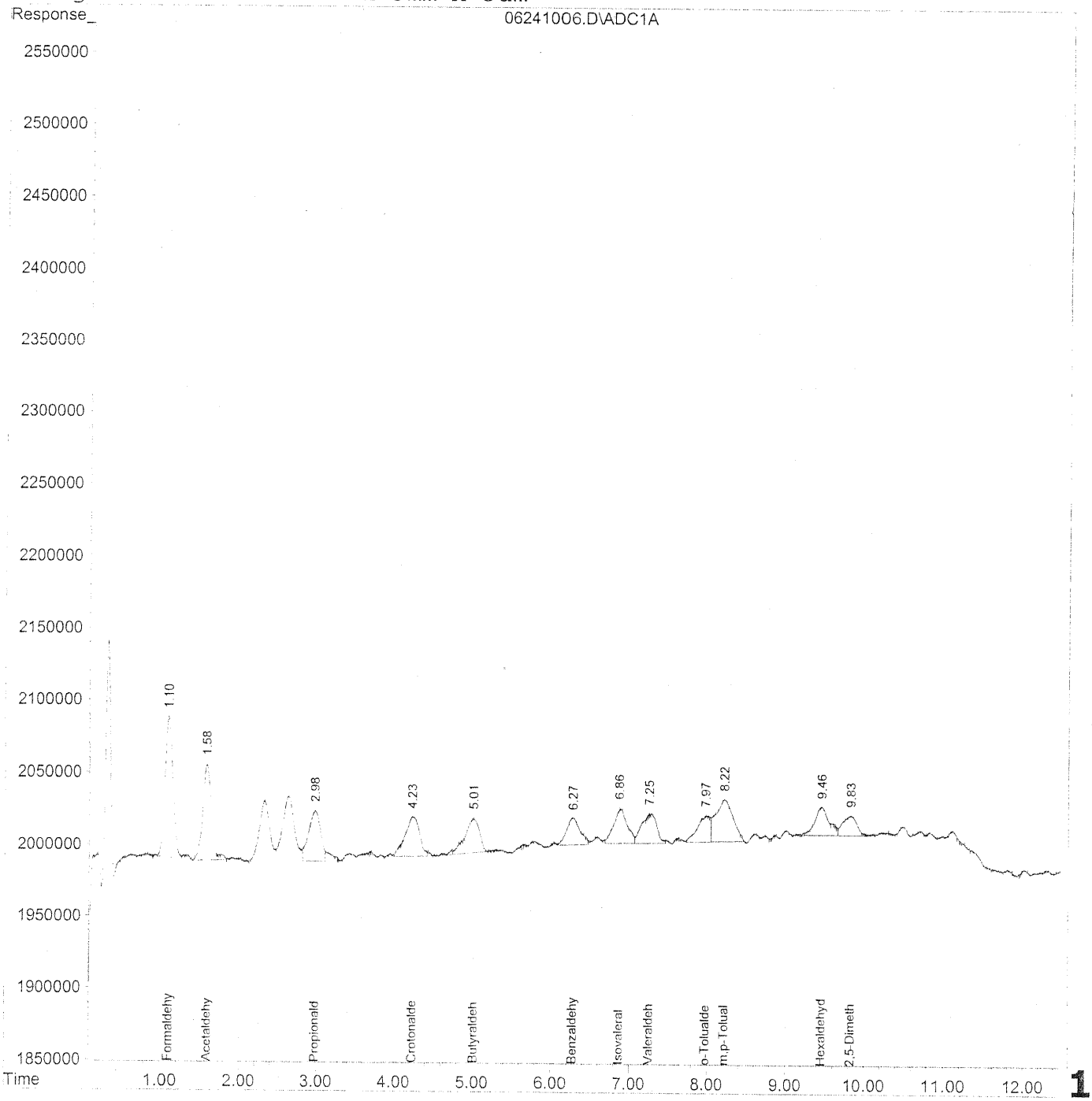
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:35 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
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\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:35 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
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

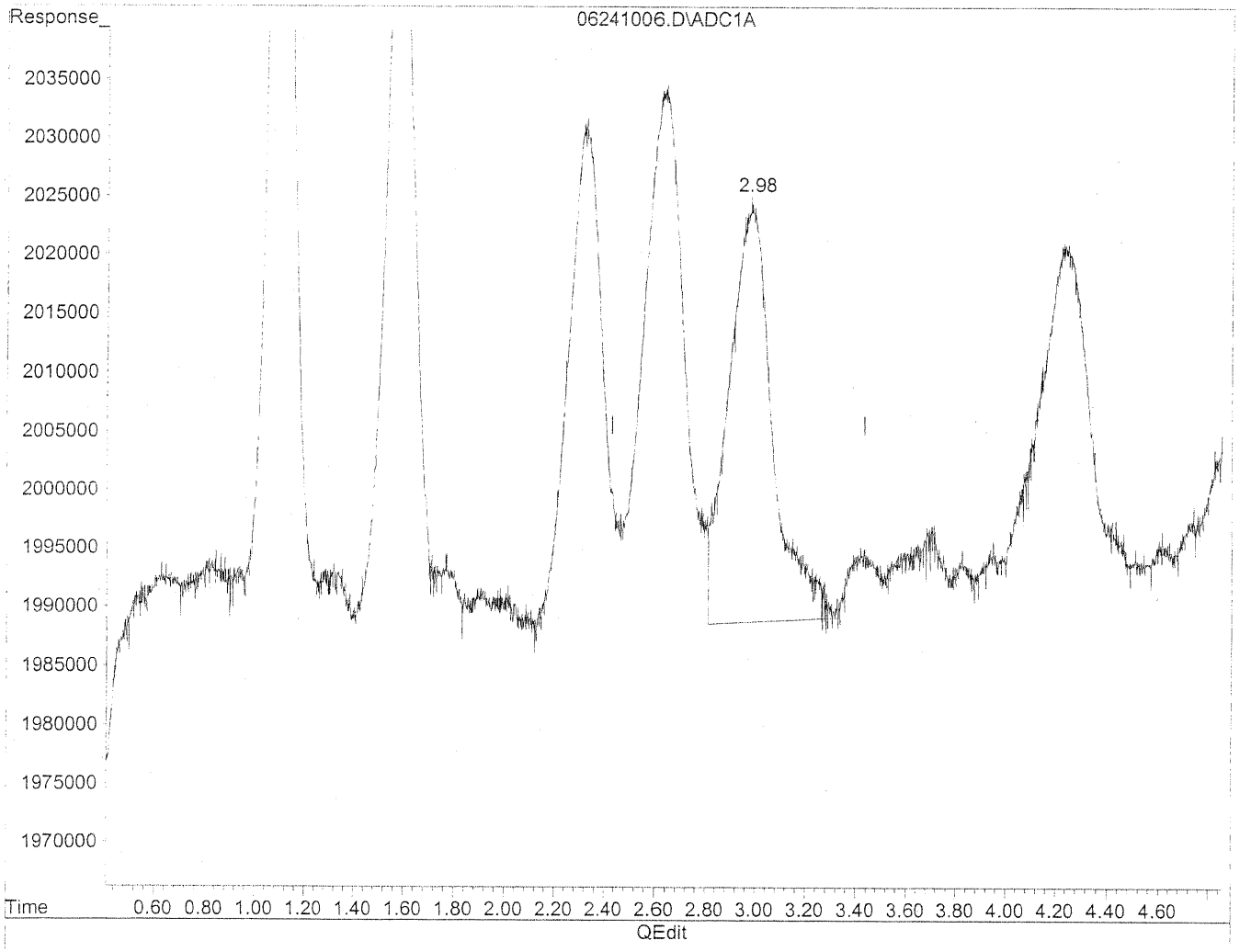
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.10	6416946	31.417	ng/ml
2)	Acetaldehyde	1.58	5390444	34.430	ng/ml
3)	Propionaldehyde	2.98	3757337	32.040	ng/mlm
4)	Crotonaldehyde	4.23	3582407	33.538	ng/mlm
5)	Butyraldehyde	5.01	3198175	32.098	ng/ml
6)	Benzaldehyde	6.27	2170695	30.083	ng/mlm
7)	Isovaleraldehyde	6.86	2928376	34.741	ng/mlm
8)	Valeraldehyde	7.25	2784817	34.841	ng/mlm
9)	o-Tolualdehyde	7.97	2066036	31.383	ng/mlm
10)	m,p-Tolualdehyde	8.22	4443968	74.153	ng/mlm
11)	Hexaldehyde	9.46	2465808	32.453	ng/mlm
12)	2,5-Dimethylbenzaldehyde	9.83	1722129	32.673	ng/mlm

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.98min 37.049ng/ml

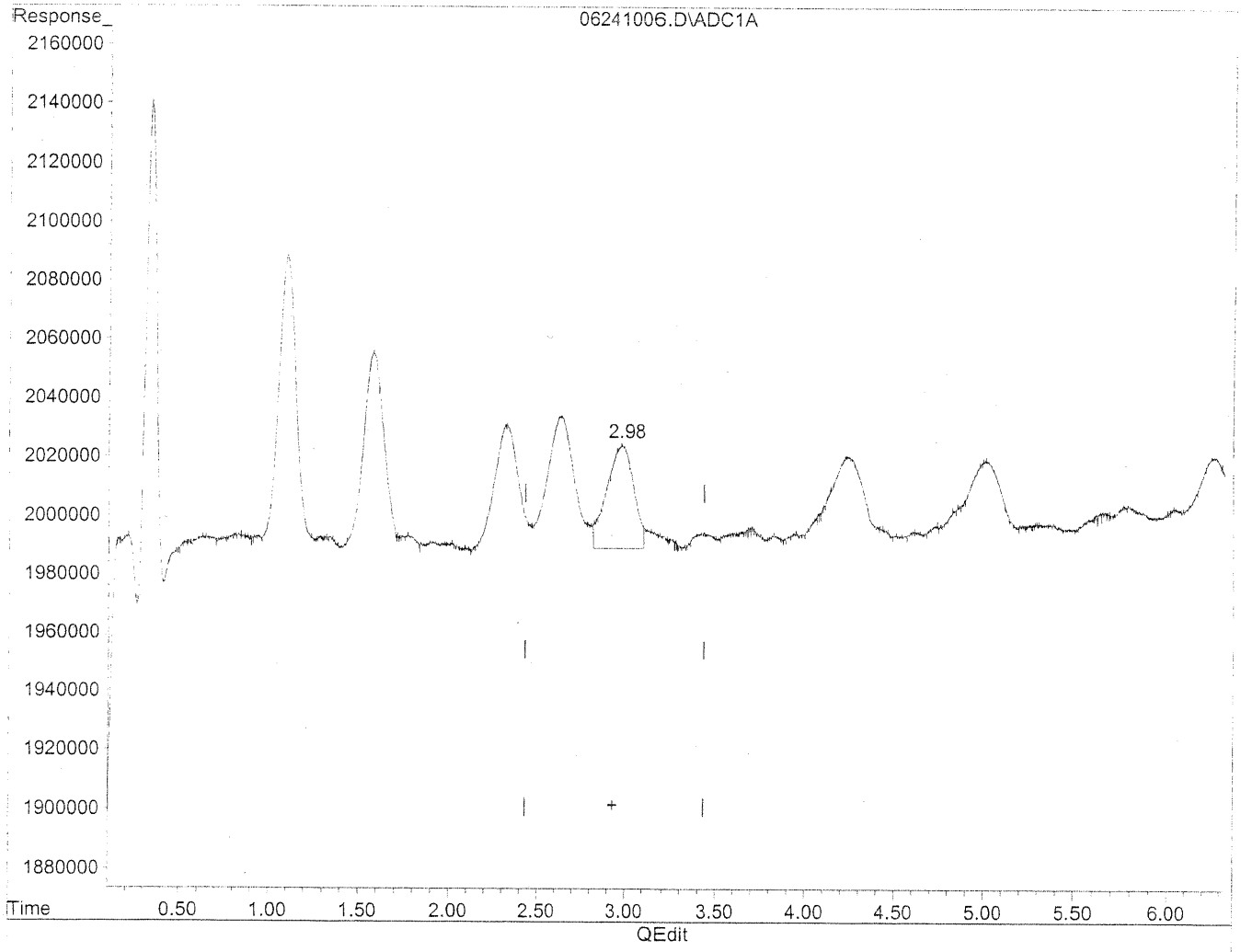
response 4344718

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(3) Propionaldehyde

2.98min 32.040ng/ml m

response 3757337

6/24/10

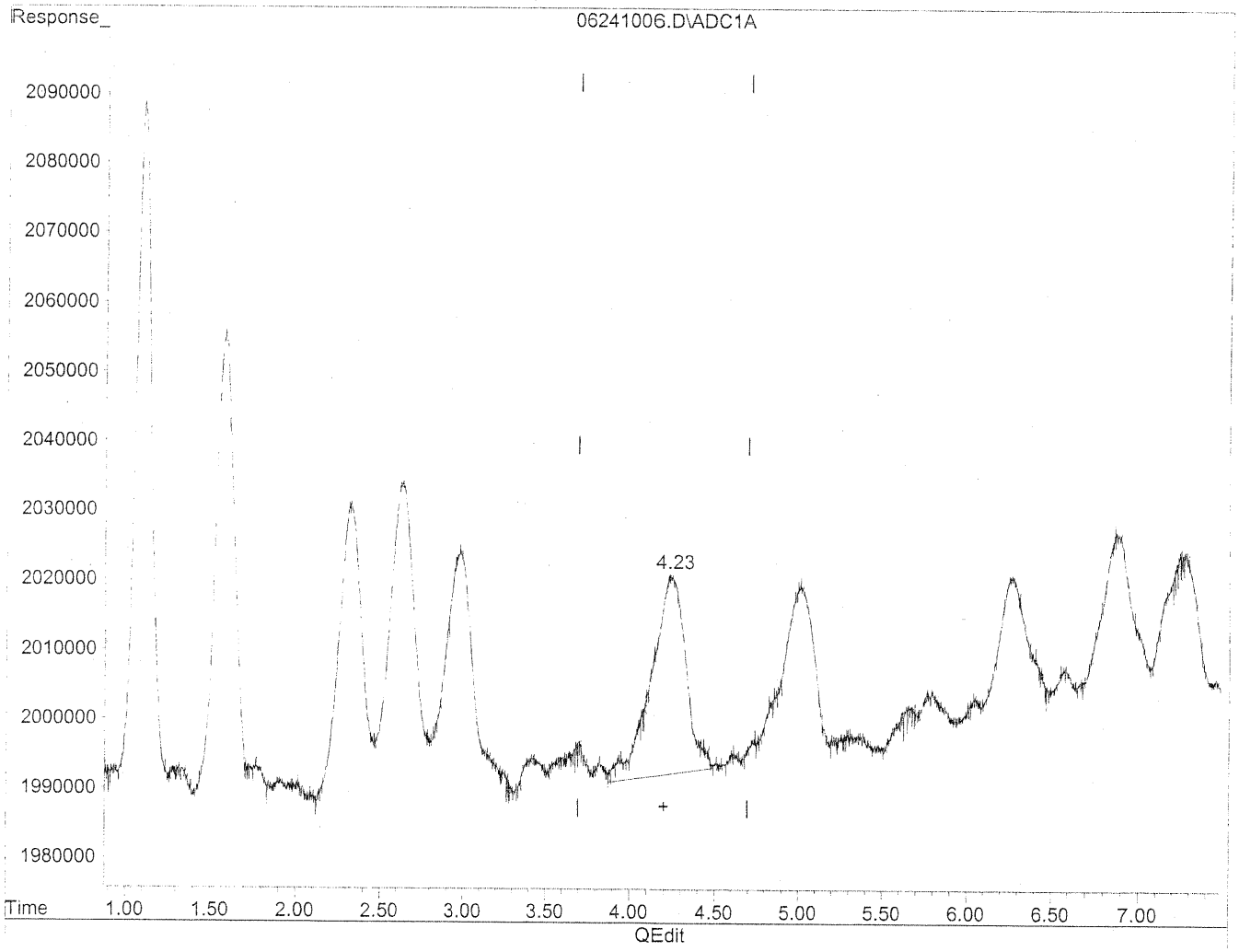
6/25/10

(+) = Expected Retention Time
06241006.D TO110610.M Thu Jun 24 13:33:49 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.24min 37.563ng/ml

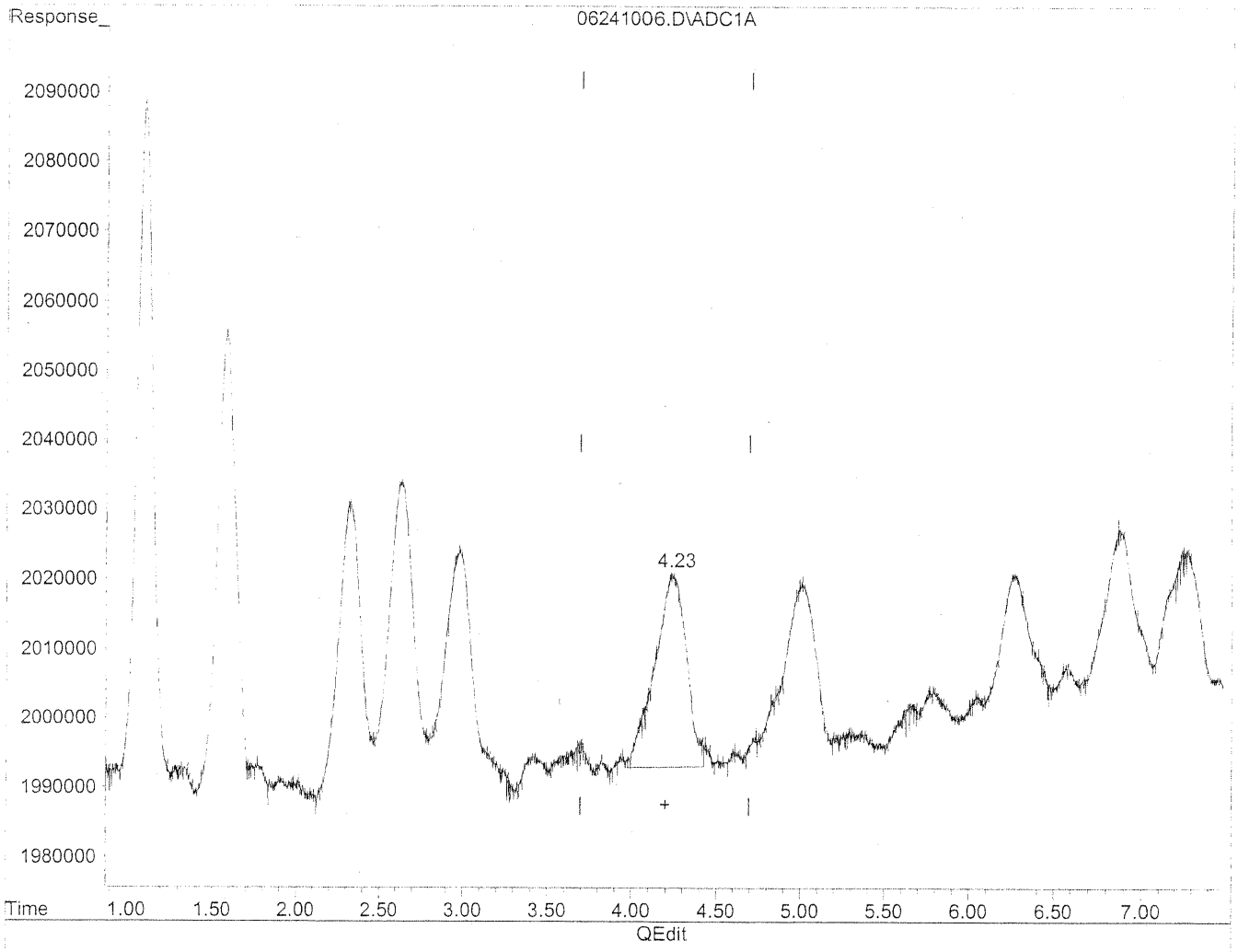
response 4012348

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.23min 33.538ng/ml m

response 3582407

tic 6/25/10

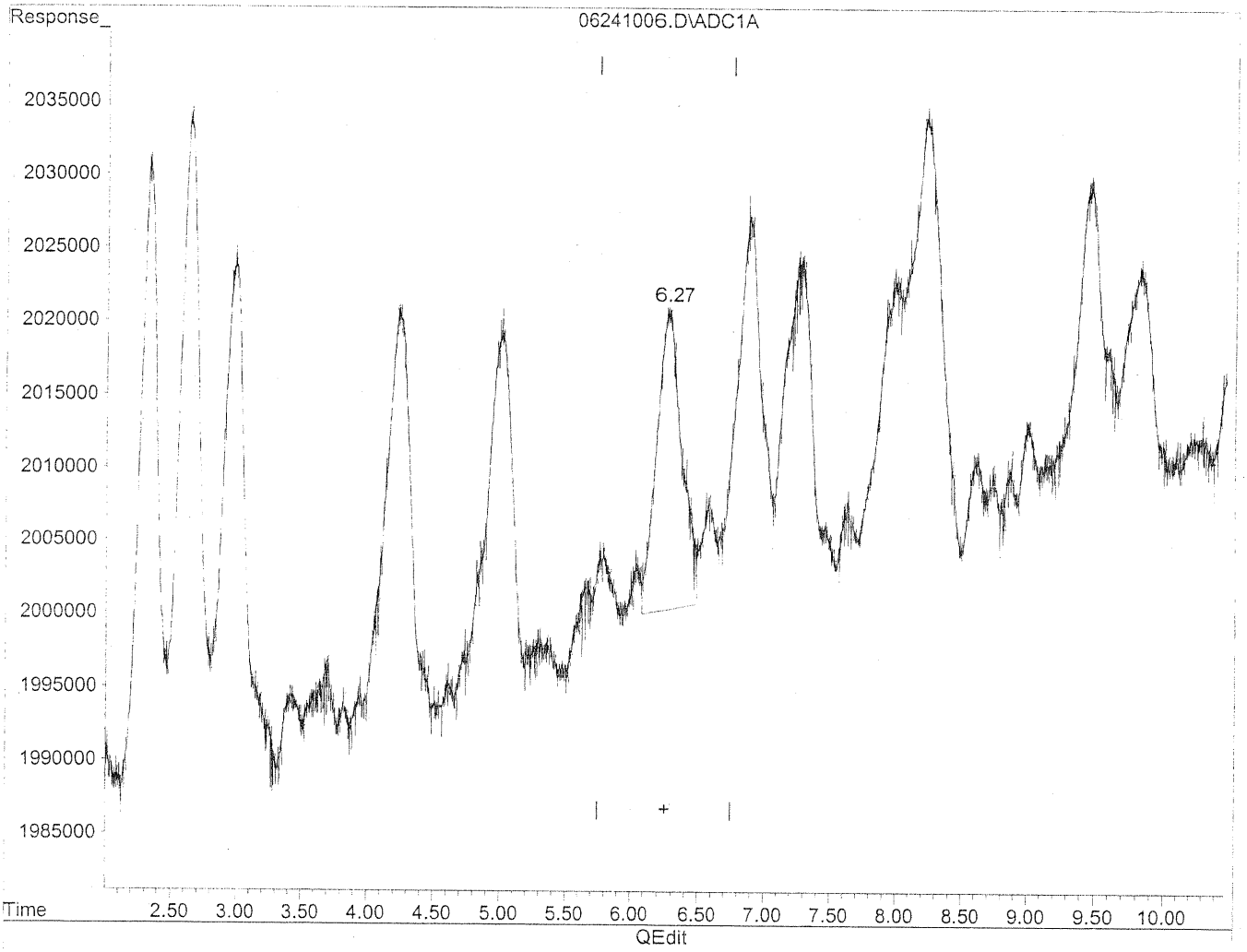
tic mp 6/25/10

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.27min 35.800ng/ml

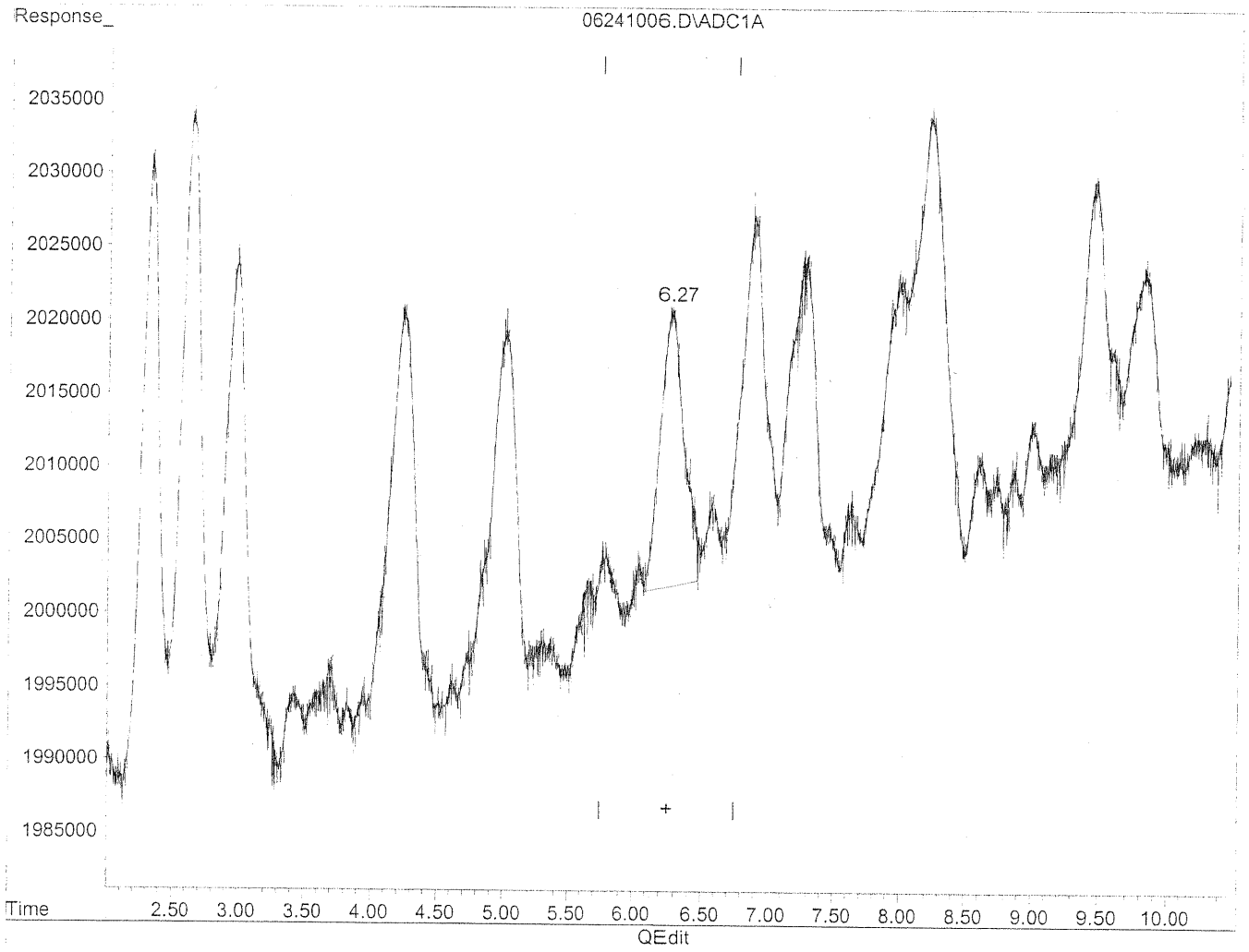
response 2583170

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.27min 30.083ng/ml m

response 2170695

4C 6/28/10

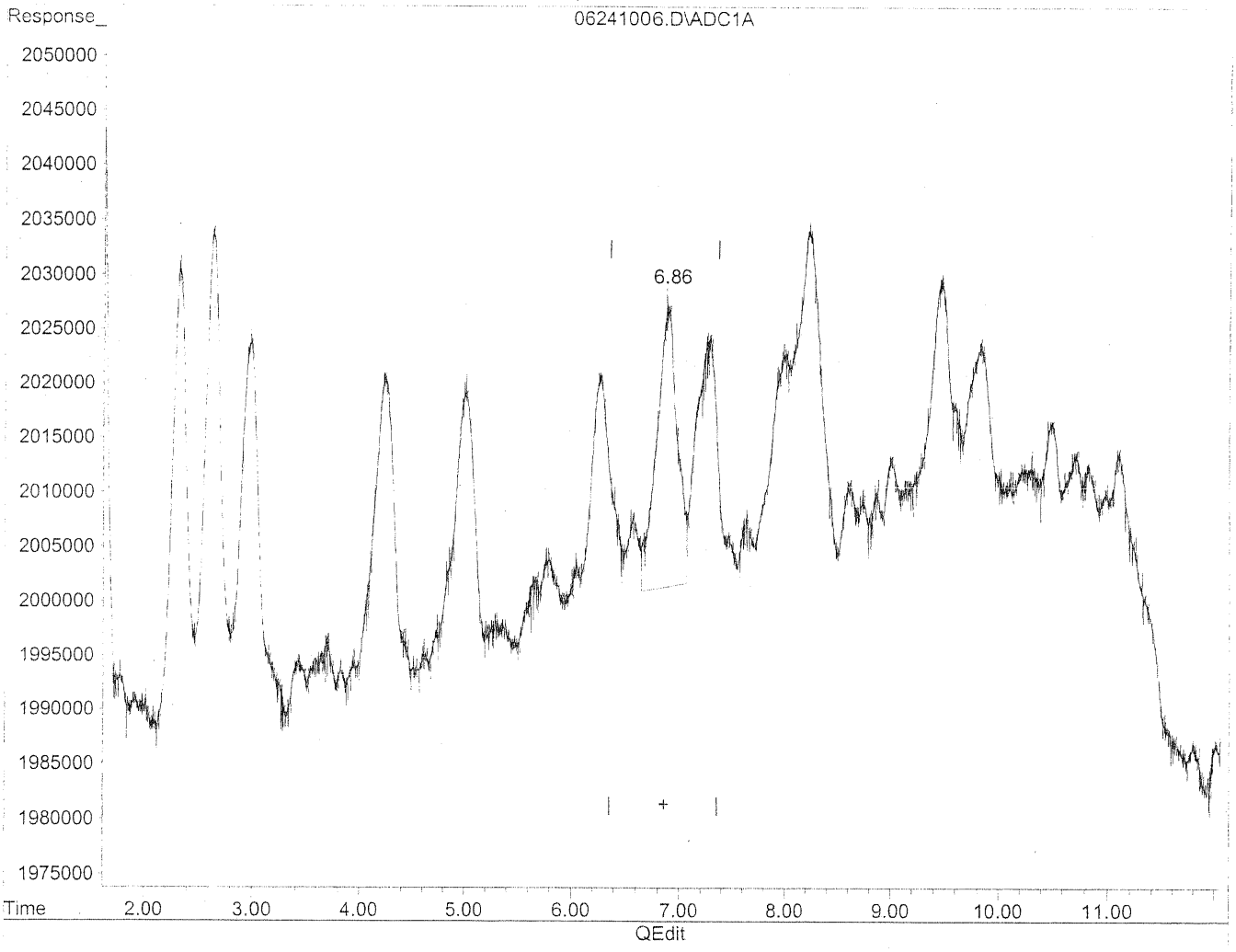
IC 6/25/10

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

6.88min 40.695ng/ml

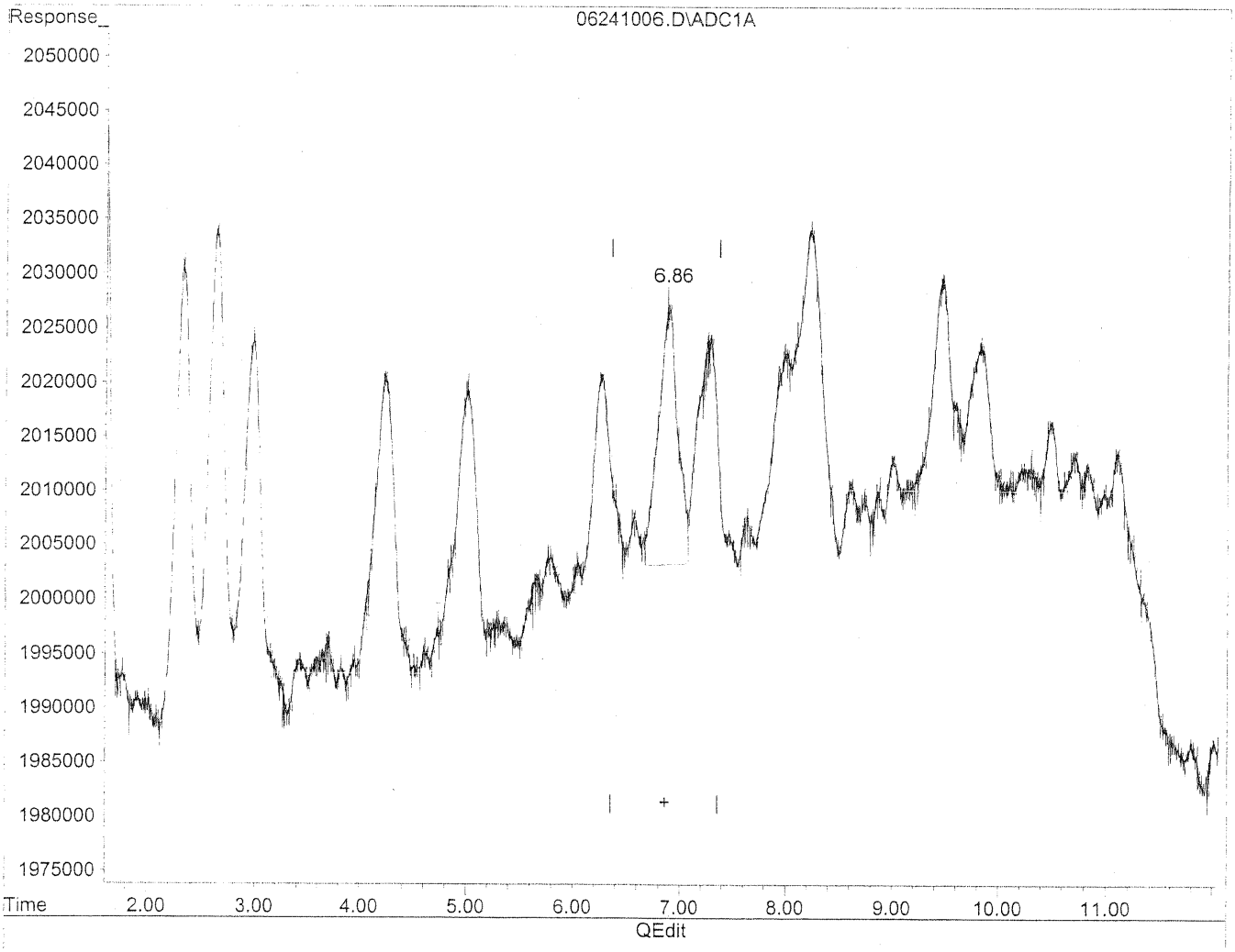
response 3430274

(+) = Expected Retention Time
06241006.D TO110610.M Thu Jun 24 13:34:23 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

6.86min 34.741ng/ml m

response 2928376

HC 6/28/10

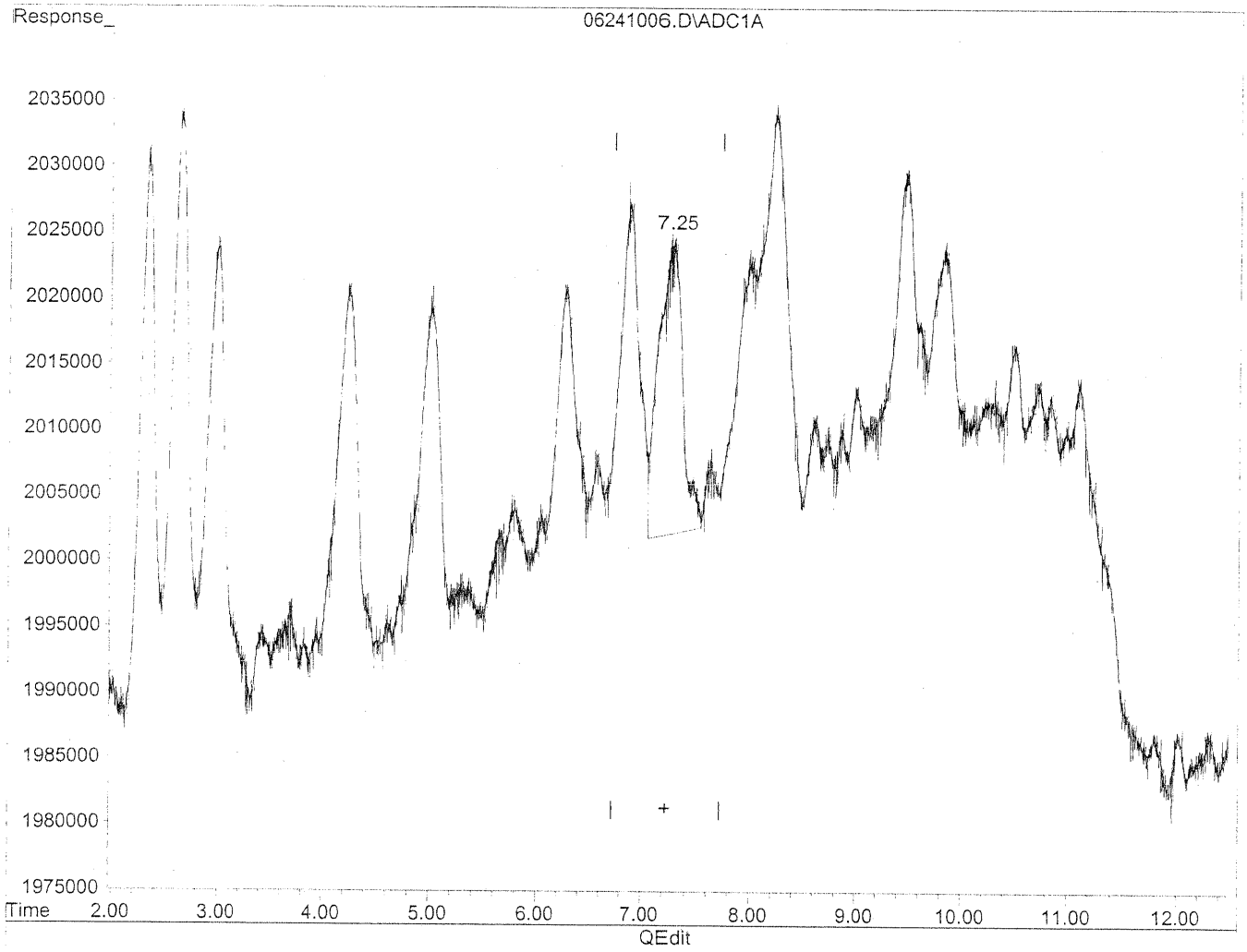
TC 6/25/10

(+) = Expected Retention Time
06241006.D TO110610.M Thu Jun 24 13:34:35 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.28min 40.180ng/ml

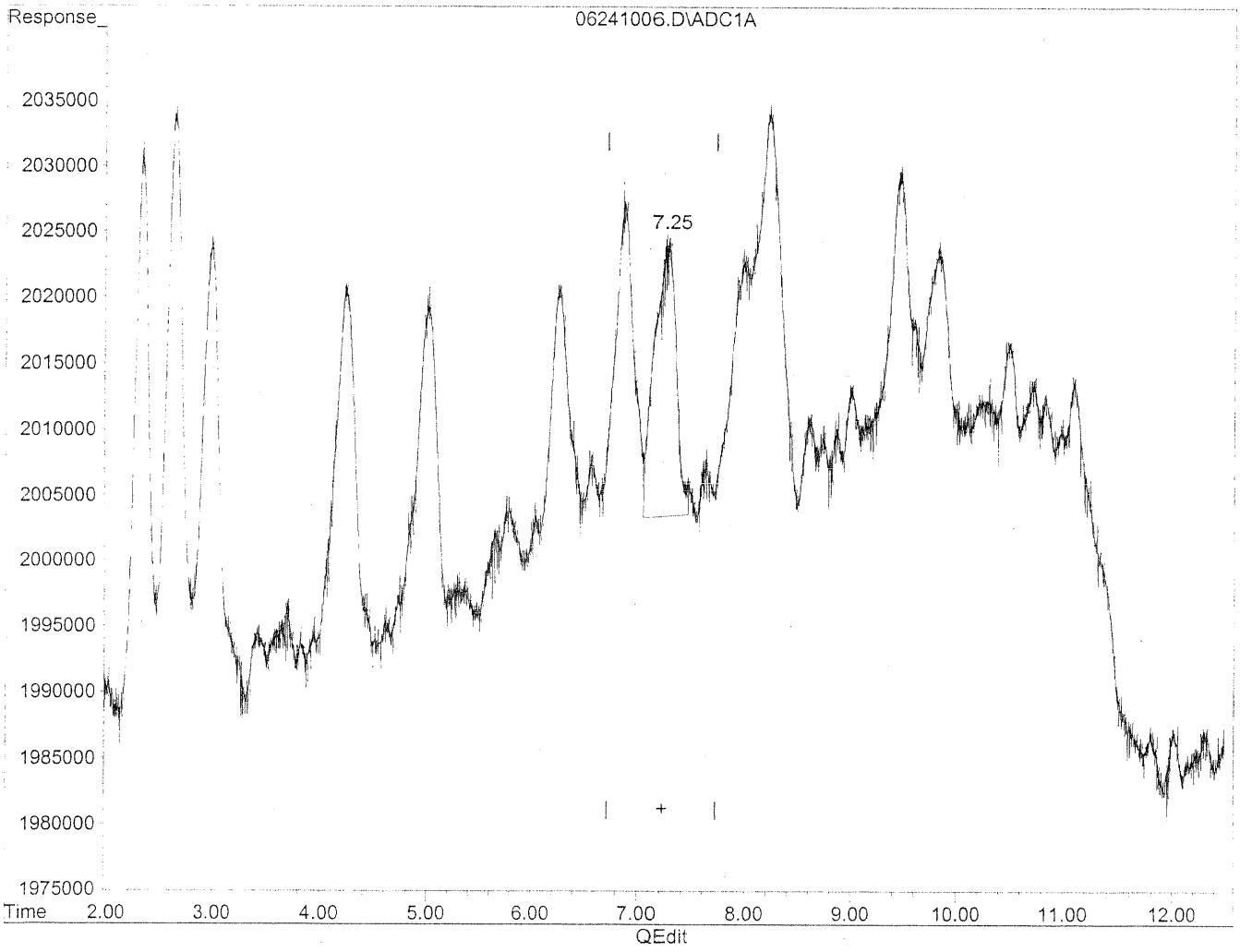
response 3211566

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.25min 34.841ng/ml m

response 2784817

6/25/10

6/25/10

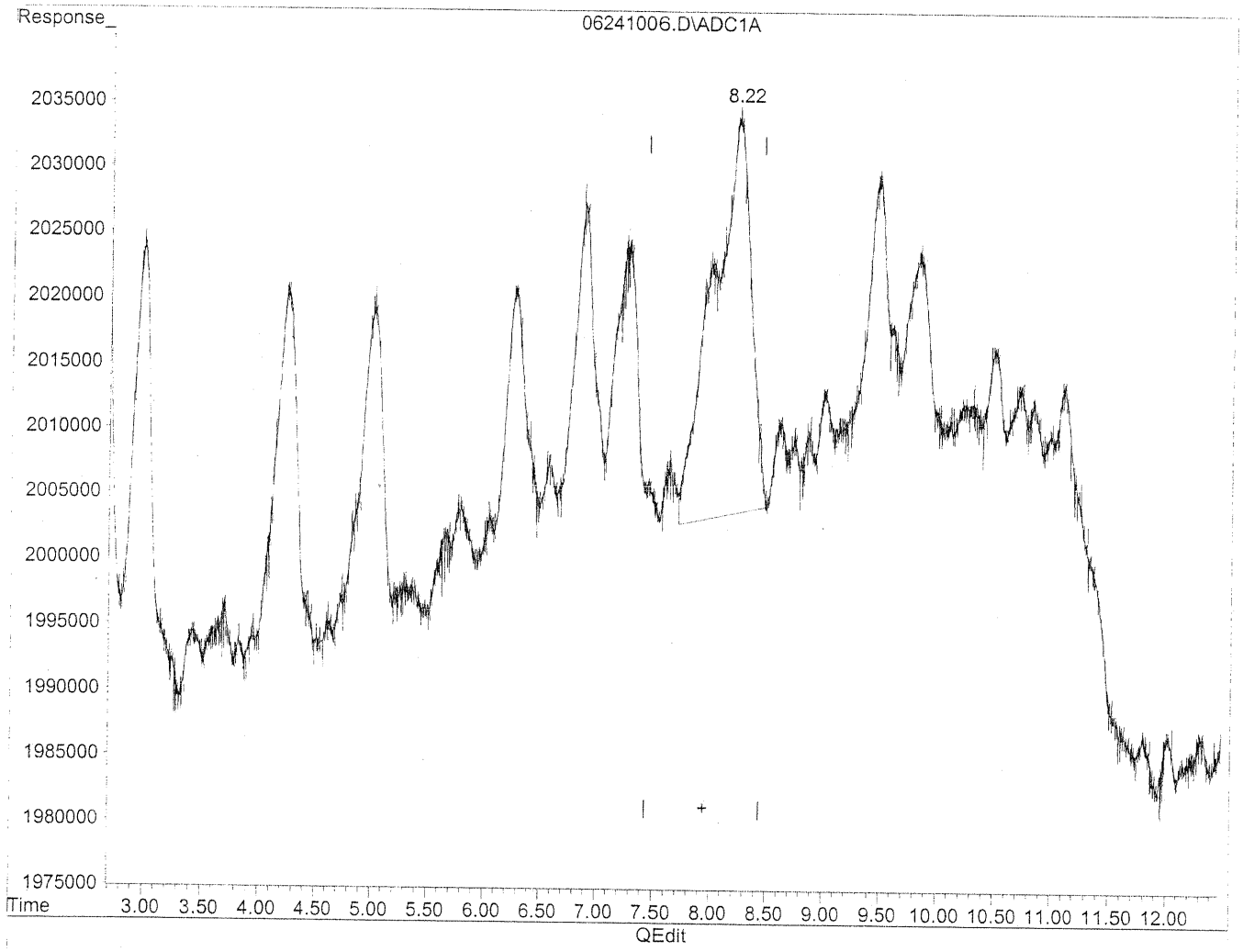
115

(+) = Expected Retention Time
06241006.D TO110610.M Thu Jun 24 13:34:52 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

8.22min 108.949ng/ml

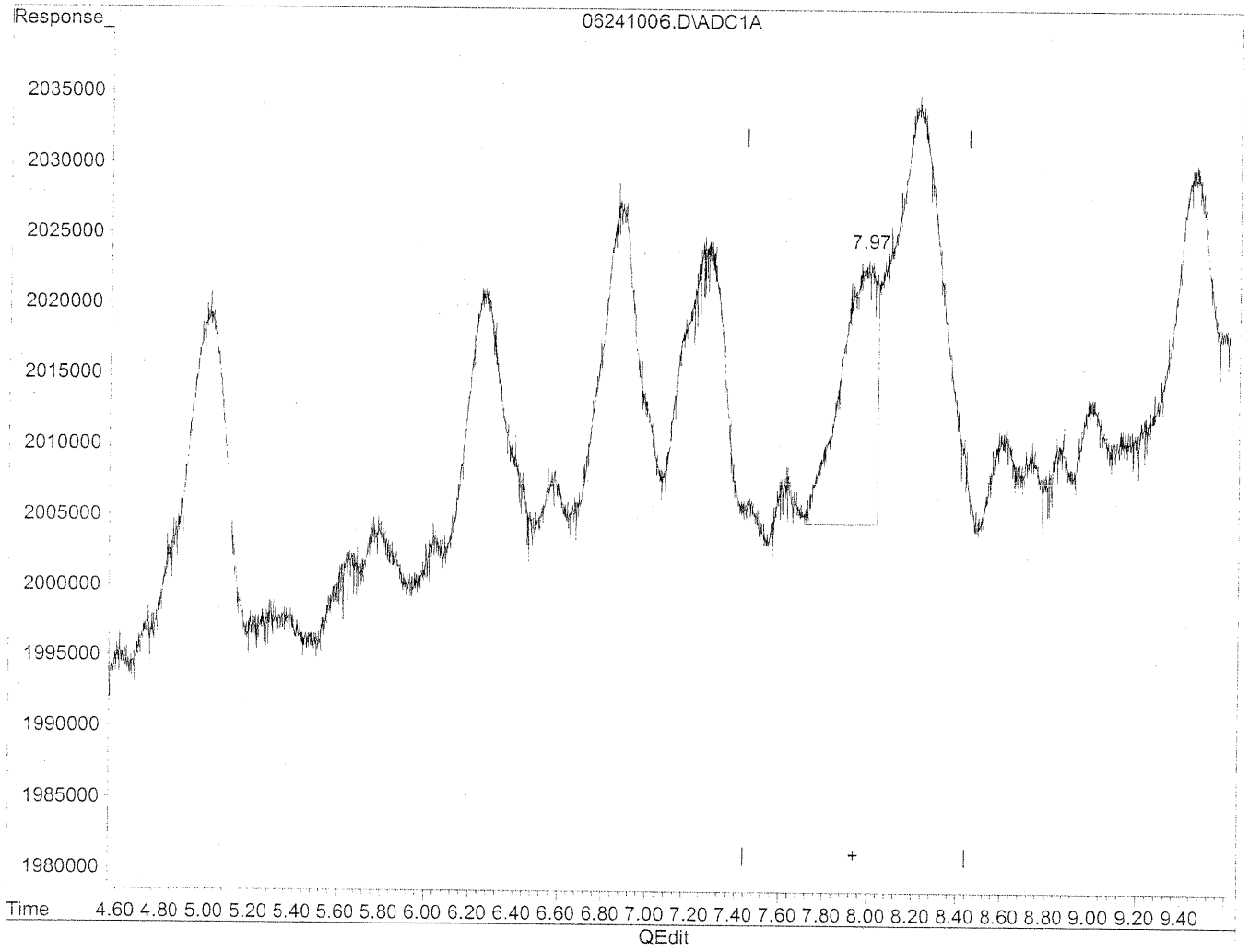
response 7172512

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

7.97min 31.383ng/ml m

response 2066036

6/25/10

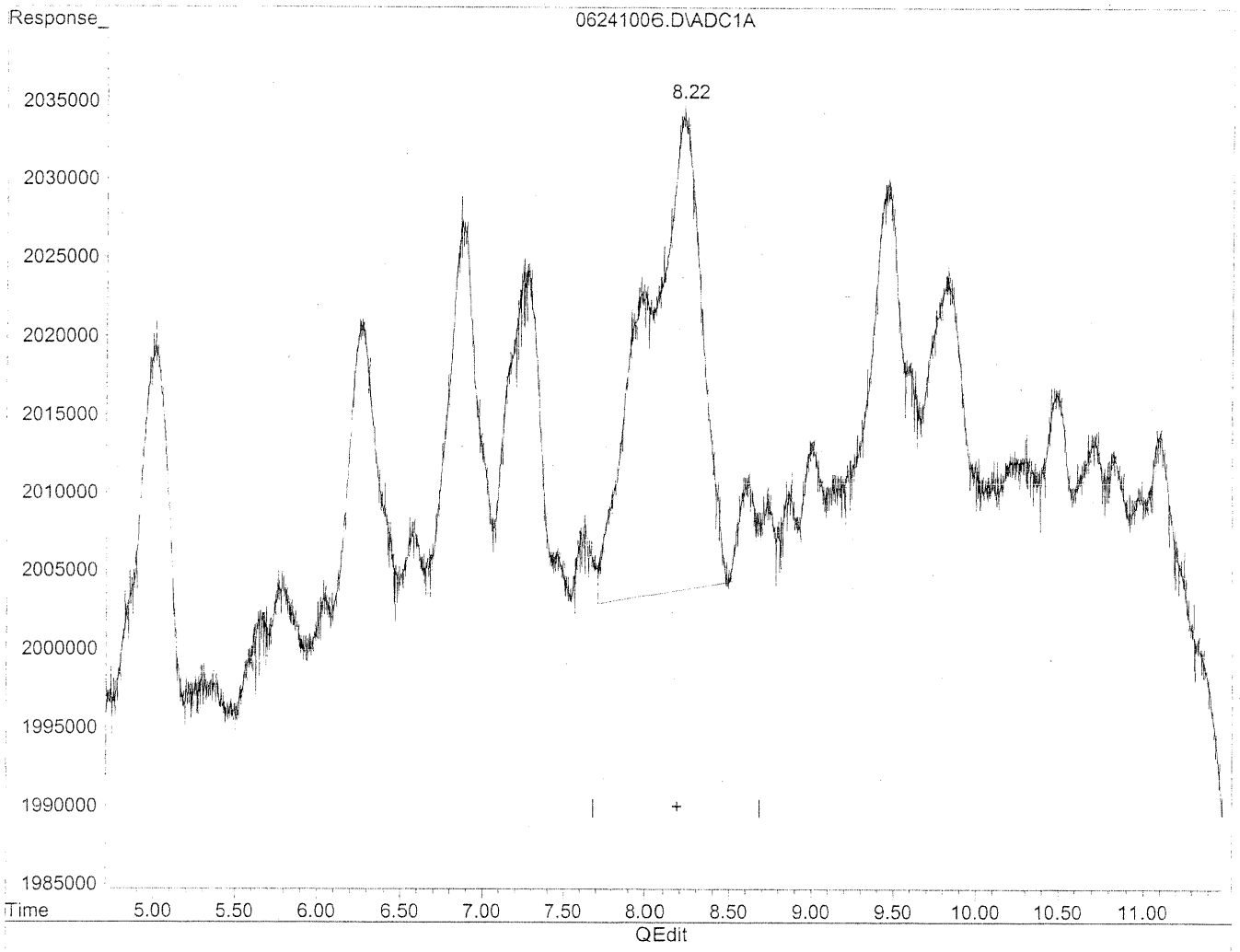
6/25/10

(+) = Expected Retention Time
06241006.D TO110610.M Thu Jun 24 13:35:09 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

8.22min 119.682ng/ml

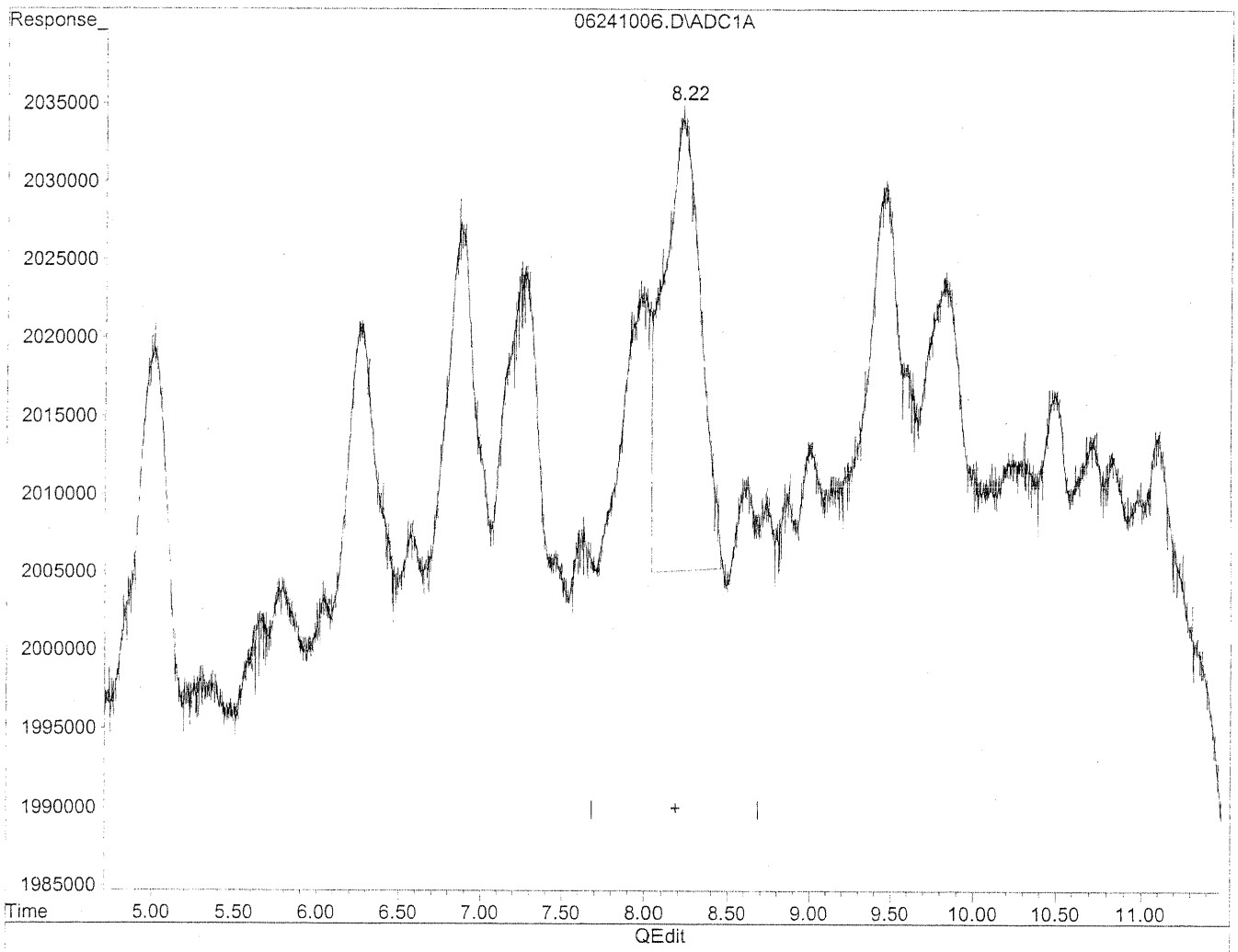
response 7172512

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

8.22min 74.153ng/ml m

response 4443968

TC 6/28/10

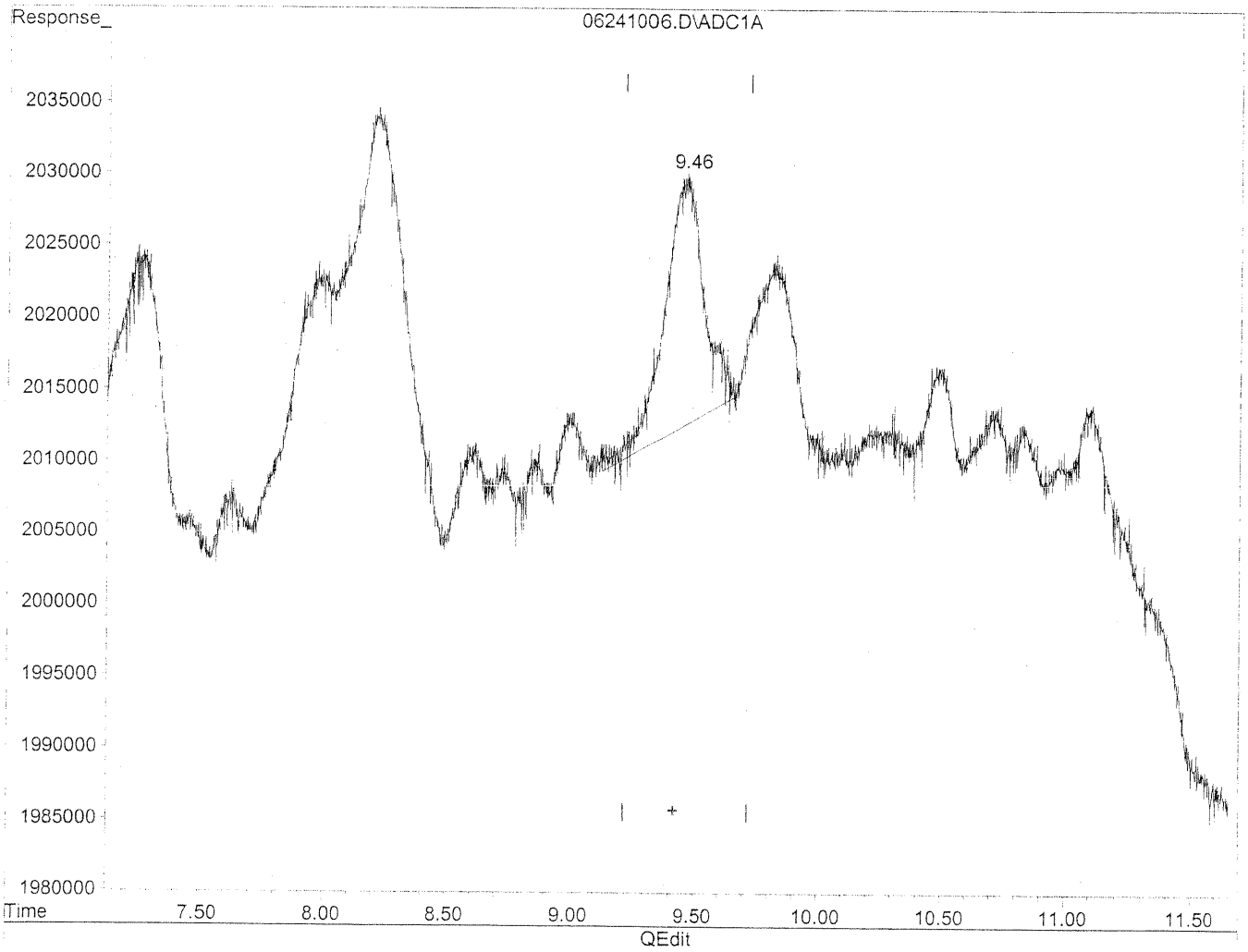
TC (MD) 6/25/10

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(11) Hexaldehyde

9.46min 24.549ng/ml

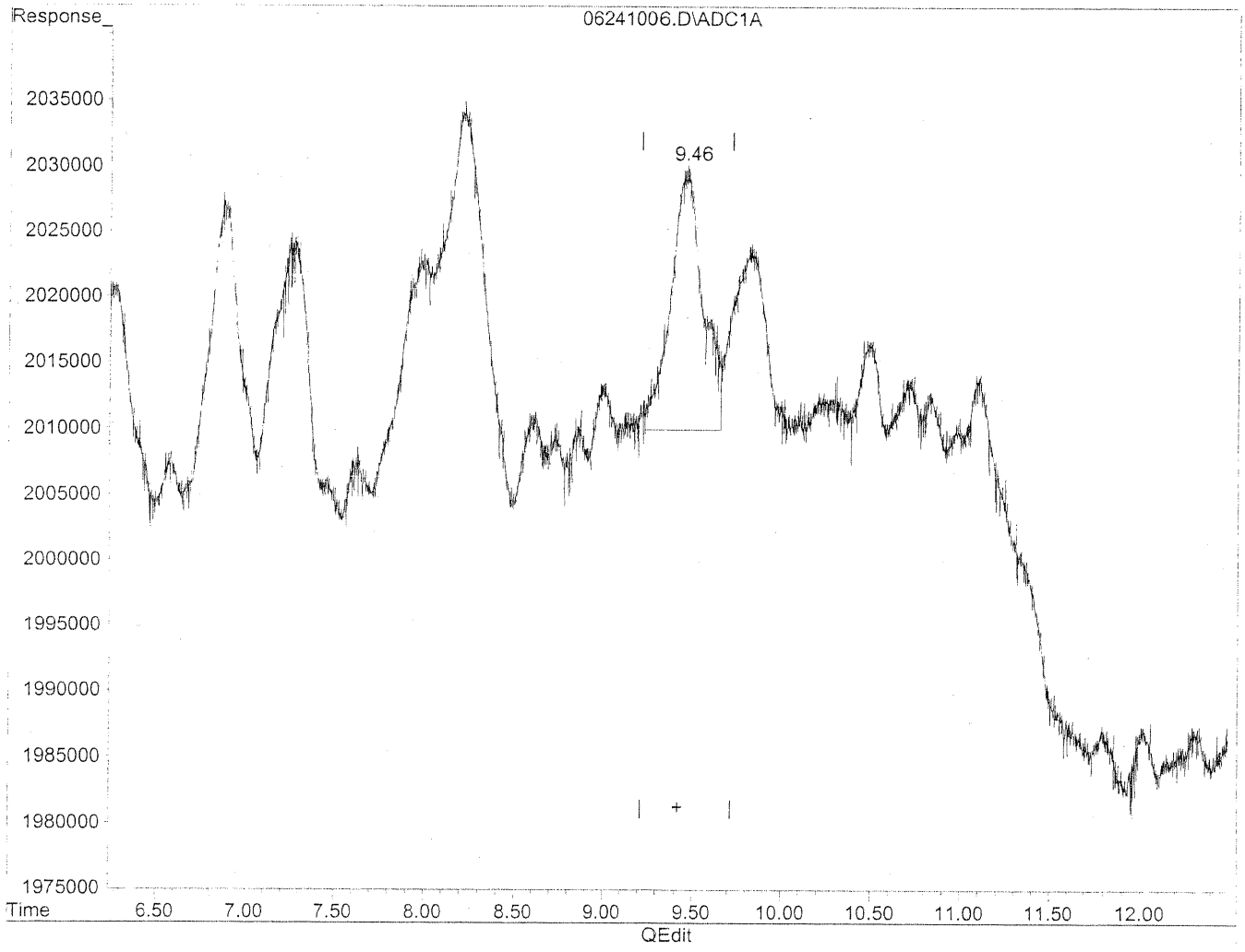
response 1865238

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(11) Hexaldehyde

9.46min 32.453ng/ml m

response 2465808

HC 6/28/10

PK 6/25/10

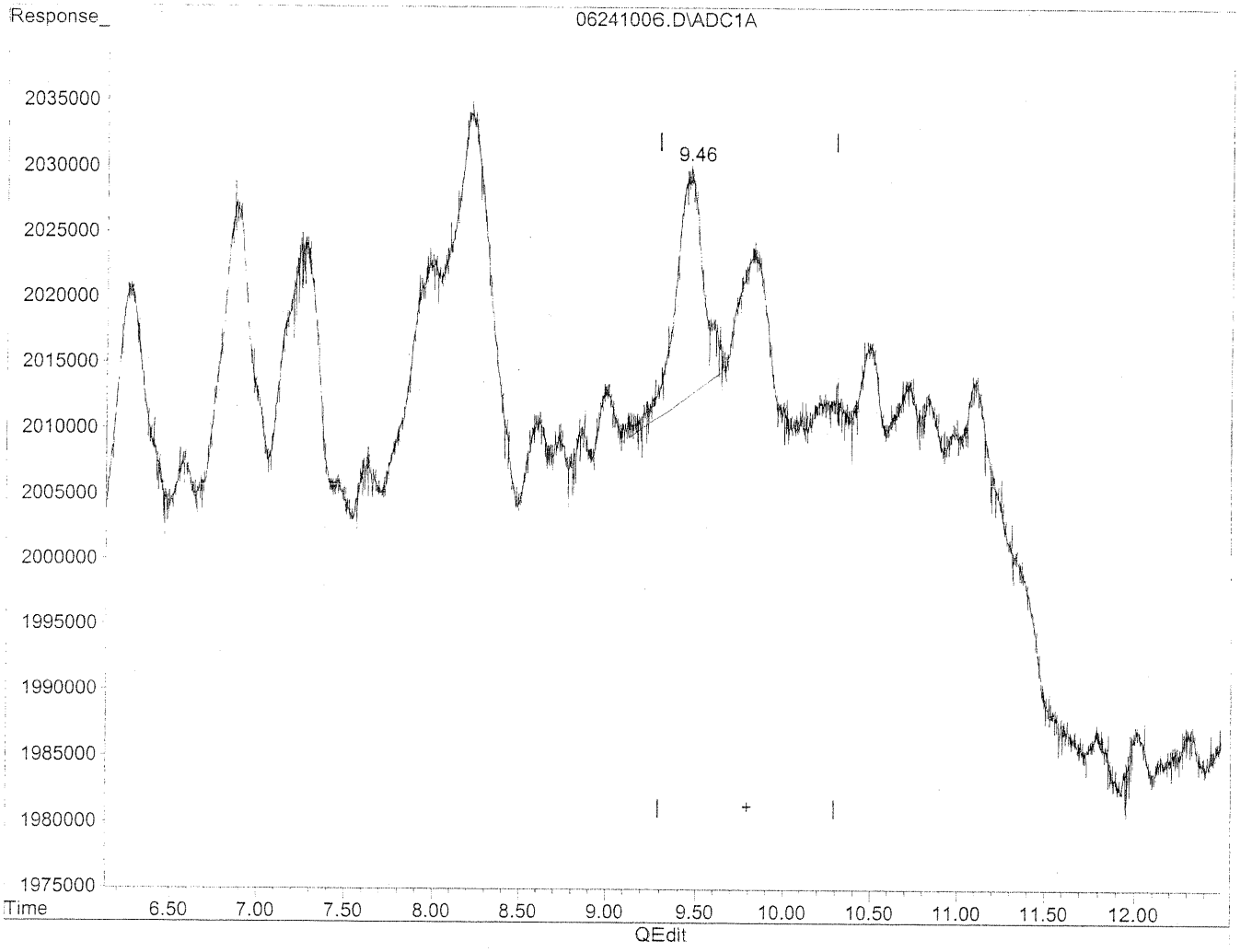
(+) = Expected Retention Time

06241006.D TO110610.M Thu Jun 24 13:35:44 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

9.46min 35.389ng/ml

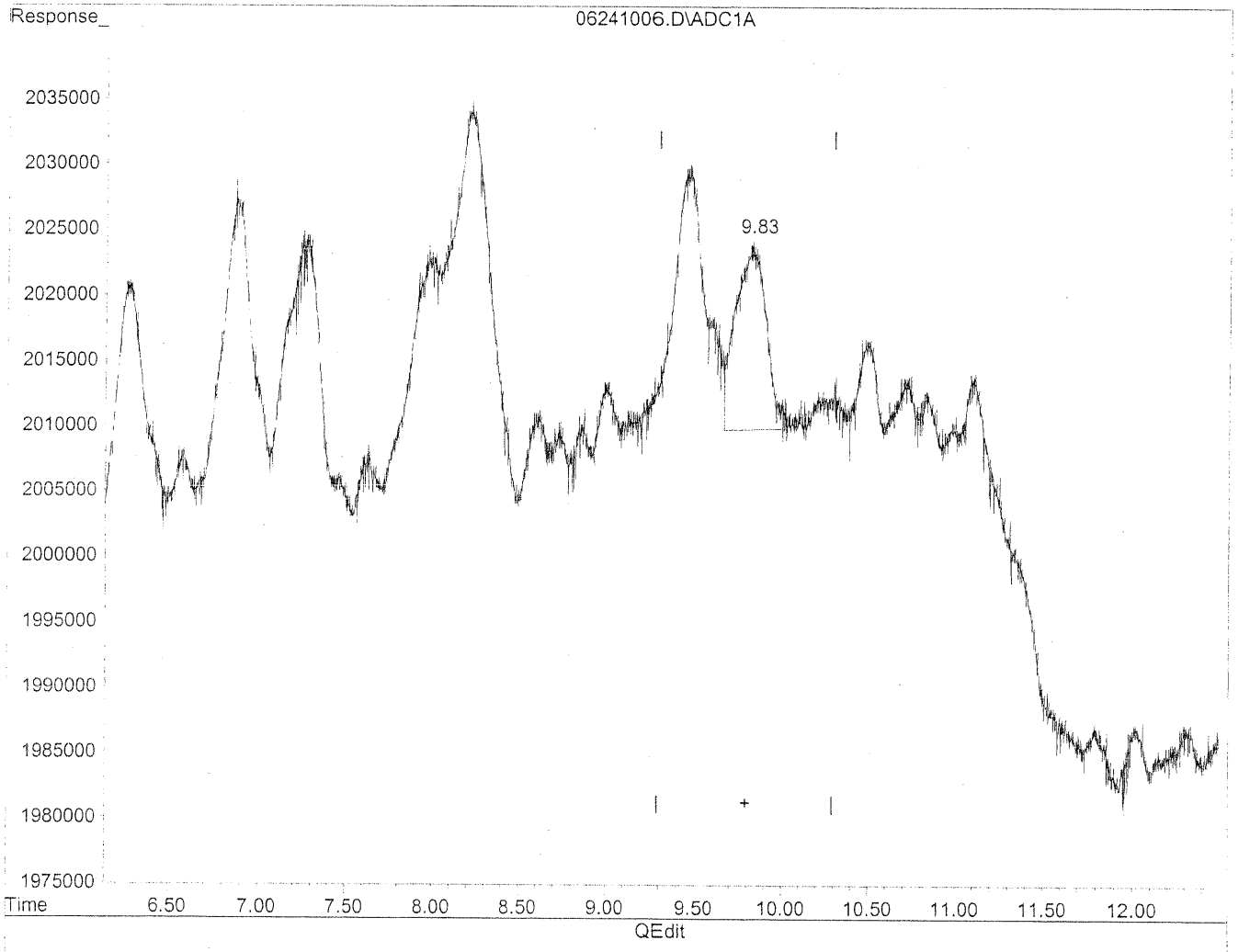
response 1865238

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241006.D Vial: 4
Acq On : 24 Jun 2010 12:31 Operator: MD
Sample : 50ng/ml TO-11A S21-03091012 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:33 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

9.83min 32.673ng/ml m

response 1722129

MP

6/25/10

(+) = Expected Retention Time

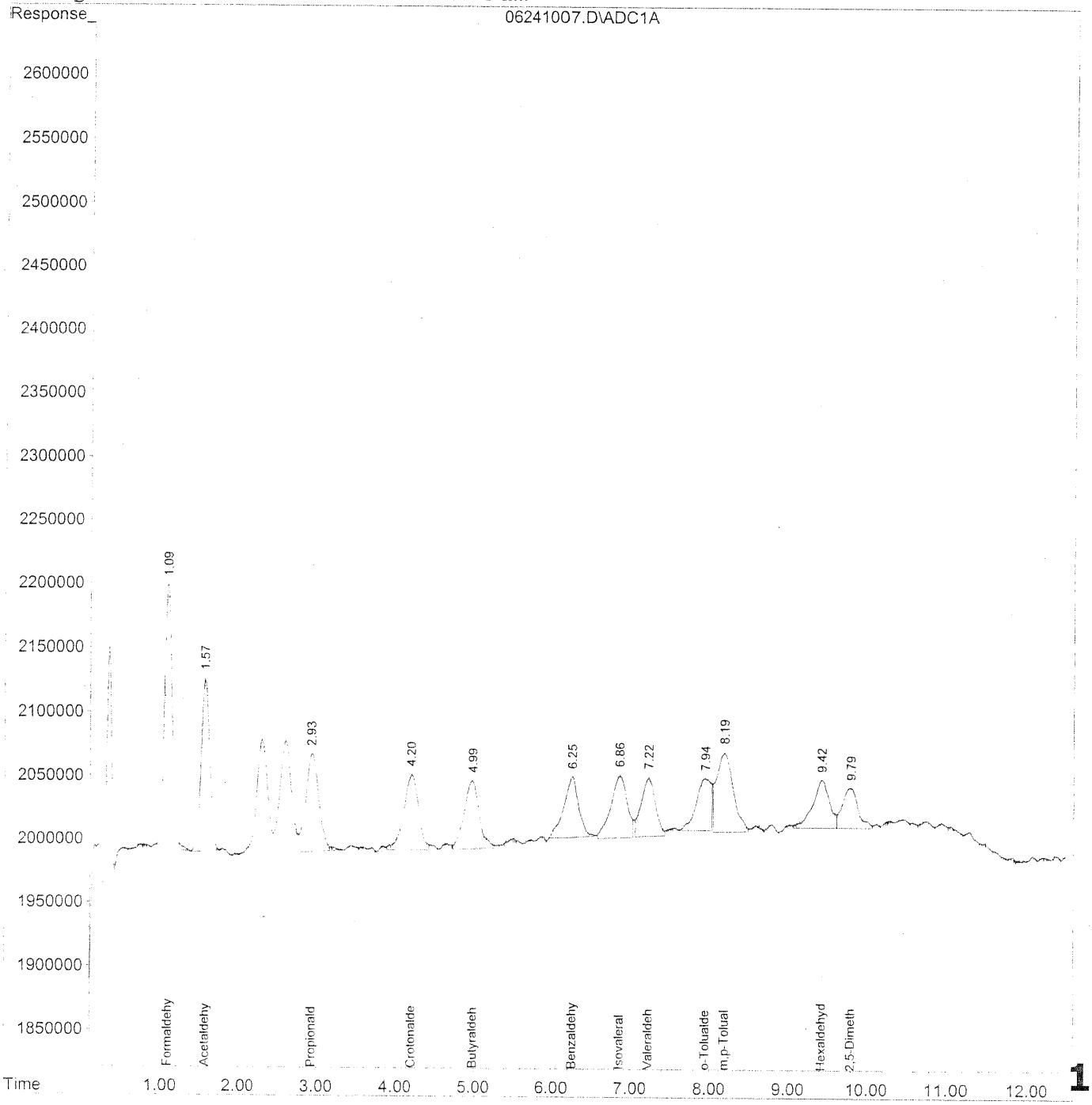
06241006.D TO110610.M Thu Jun 24 13:35:56 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:49 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
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\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:49 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
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

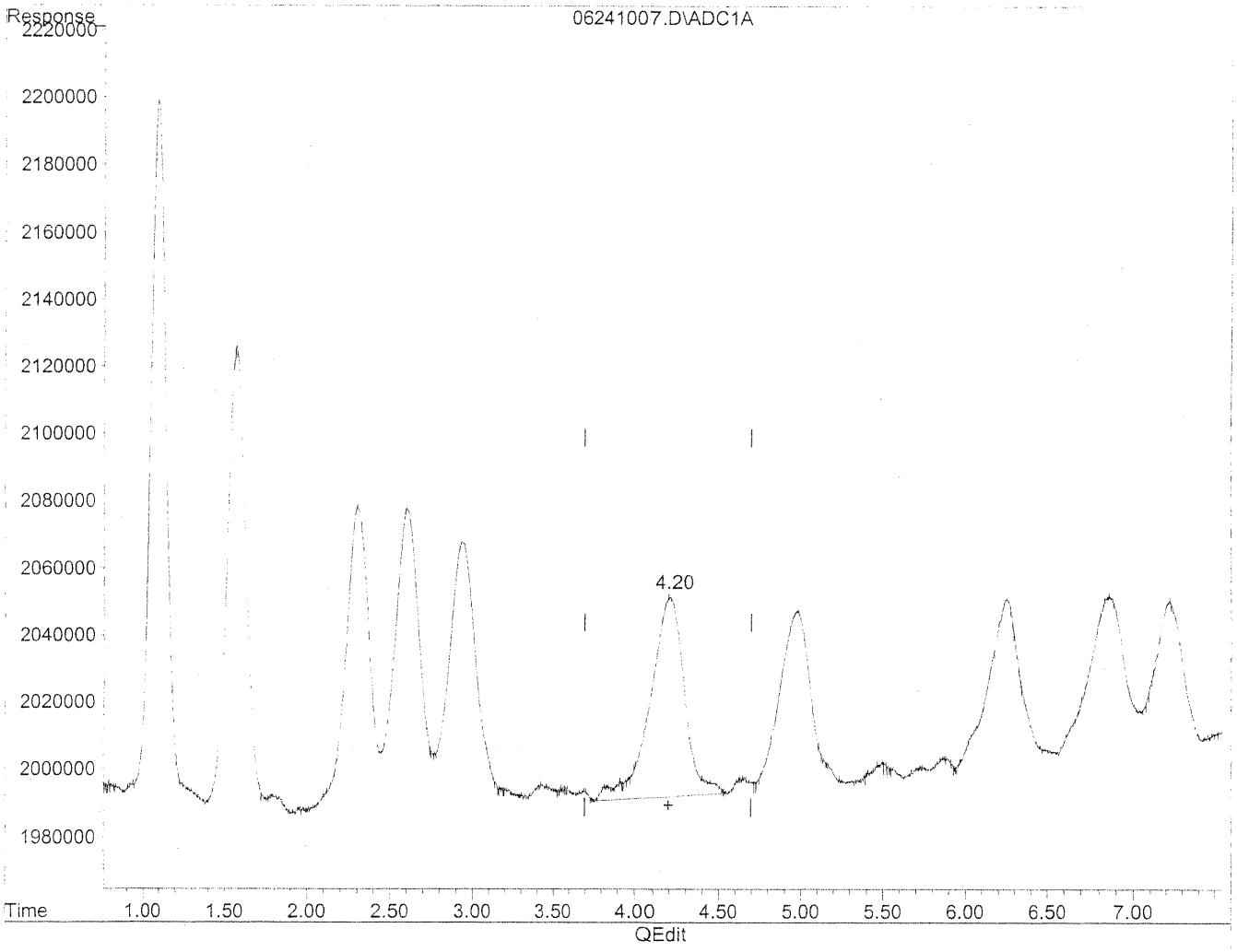
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.09	13684484	67.095	ng/ml
2)	Acetaldehyde	1.57	10561892	67.389	ng/ml
3)	Propionaldehyde	2.94	8914731	75.959	ng/ml
4)	Crotonaldehyde	4.20	7957674	74.443	ng/mlm
5)	Butyraldehyde	4.99	7348500	74.074	ng/mlm
6)	Benzaldehyde	6.25	6214018	86.389	ng/mlm
7)	Isovaleraldehyde	6.86	7015814	82.866	ng/mlm
8)	Valeraldehyde	7.22	5776234	72.167	ng/mlm
9)	o-Tolualdehyde	7.94	5309971	81.051	ng/mlm
10)	m,p-Tolualdehyde	8.19	9000241	149.541	ng/mlm
11)	Hexaldehyde	9.42	5629872	73.979	ng/mlm
12)	2,5-Dimethylbenzaldehyde	9.79	4497151	85.755	ng/mlm

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:36 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.20min 79.259ng/ml

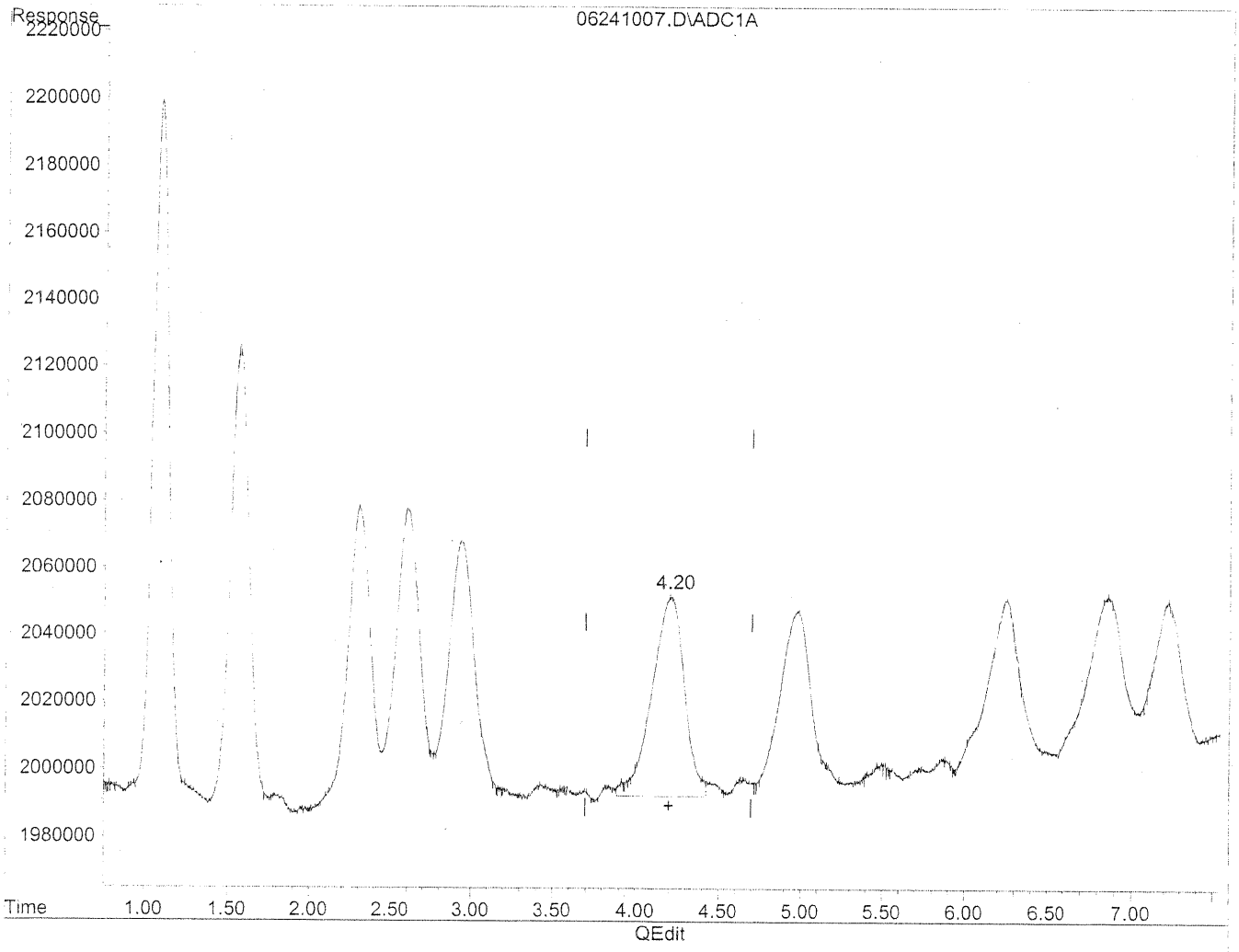
response 8472541

(+) = Expected Retention Time
06241007.D TO110610.M Thu Jun 24 13:36:51 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:36 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.20min 74.443ng/ml m

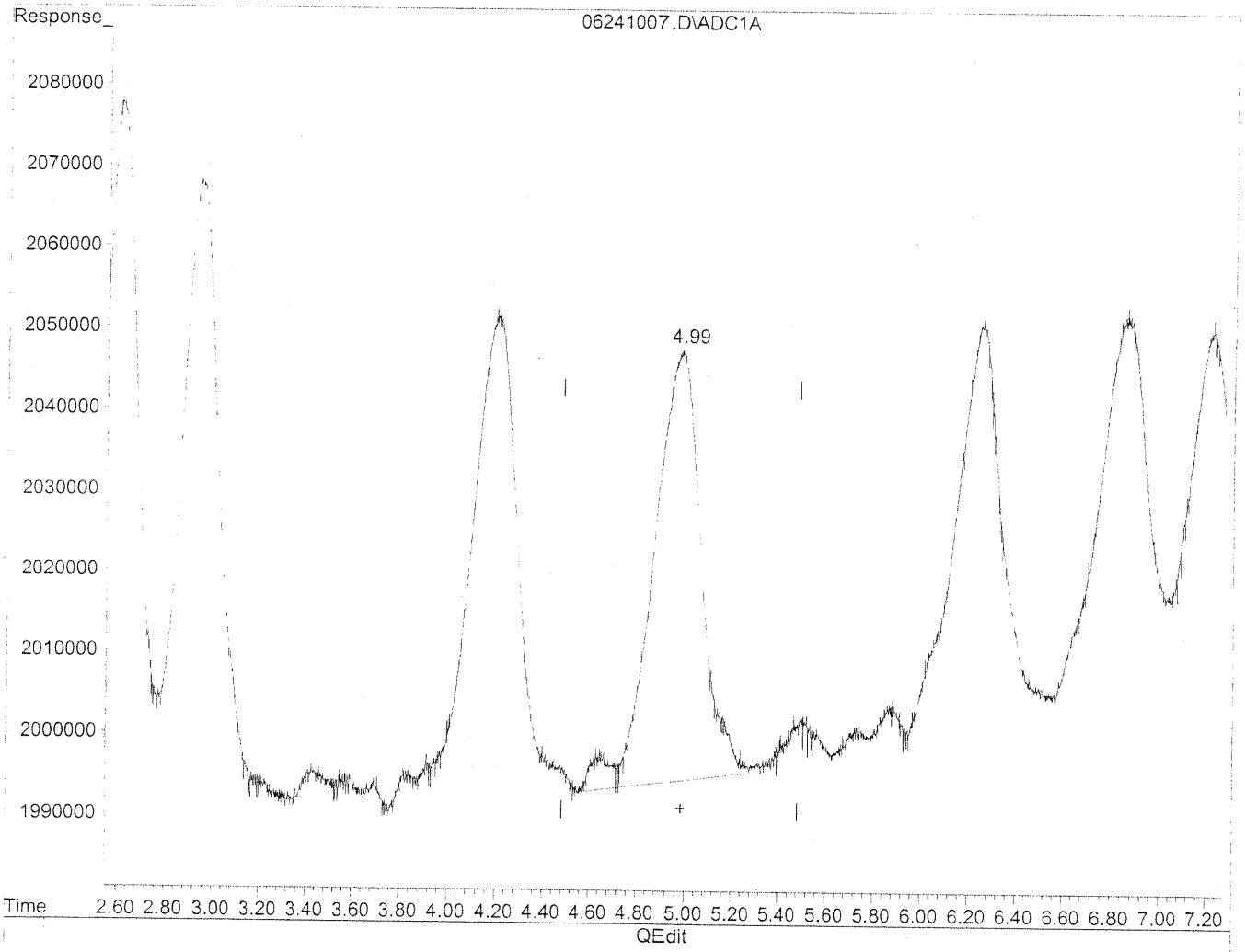
response 7957674

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:36 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.98min 73.495ng/ml

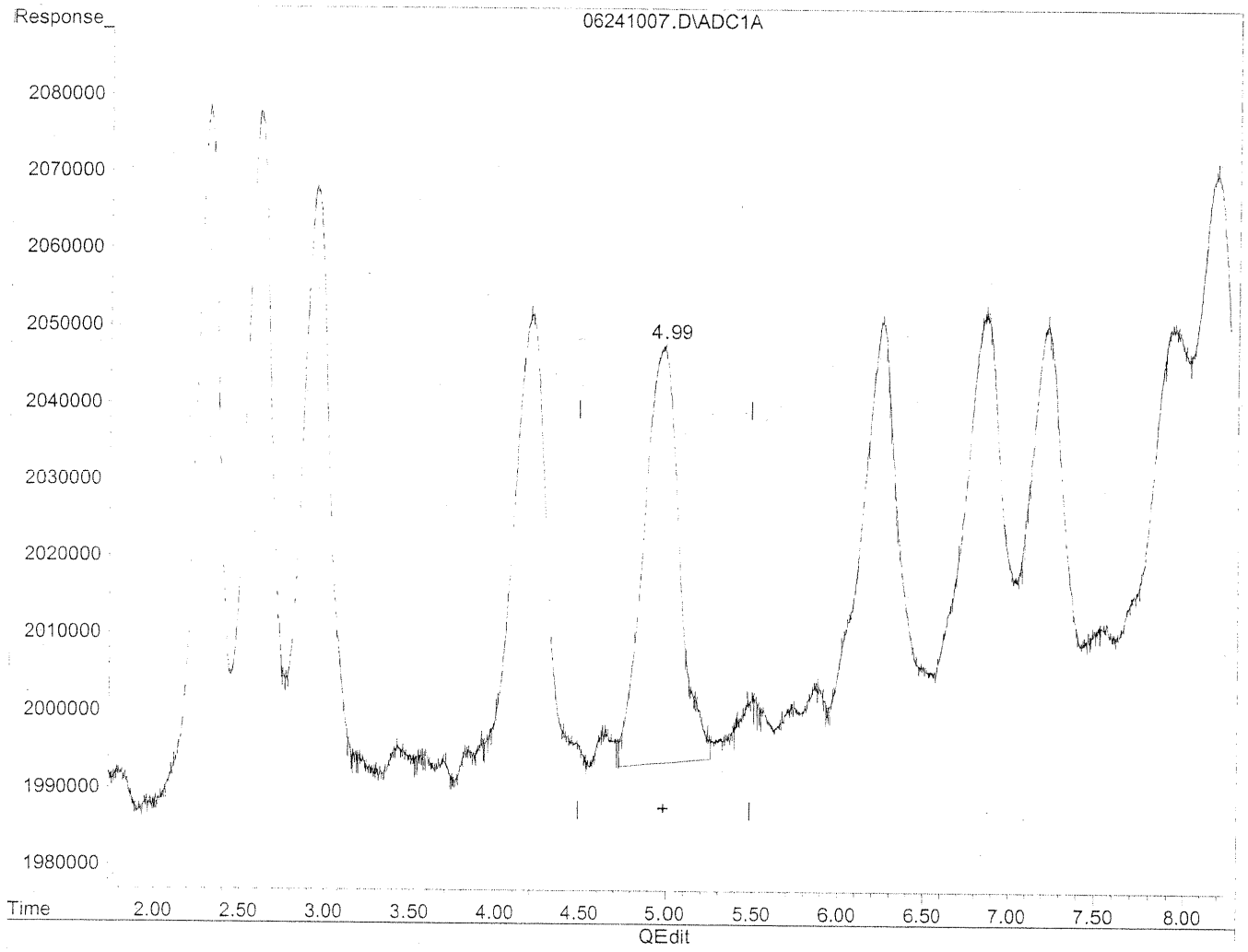
response 7291083

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:39 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.99min 74.074ng/ml m

response 7348500

6/28/10 6/25/10

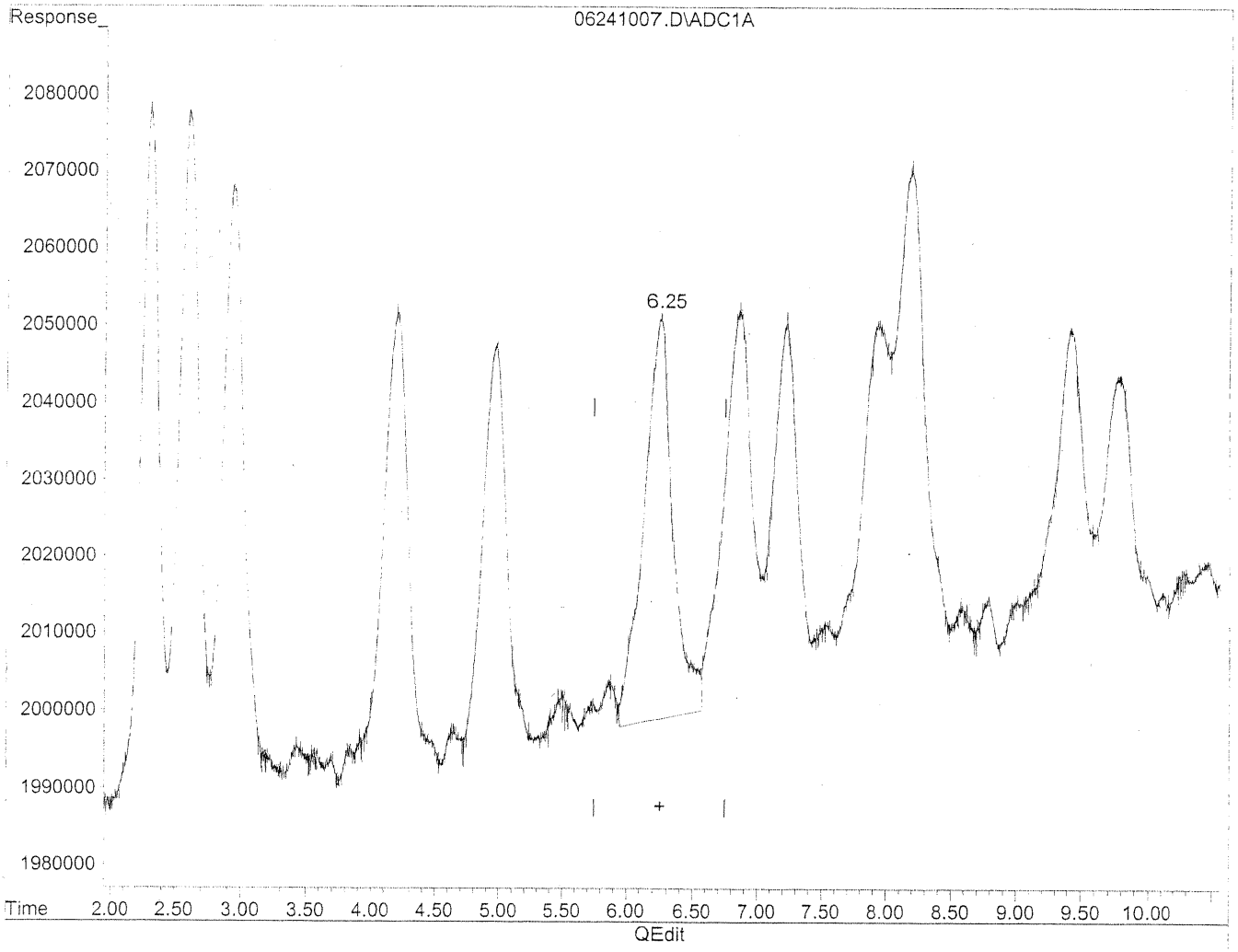
(+) = Expected Retention Time

06241007.D TO110610.M Thu Jun 24 16:27:57 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:36 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.25min 108.048ng/ml

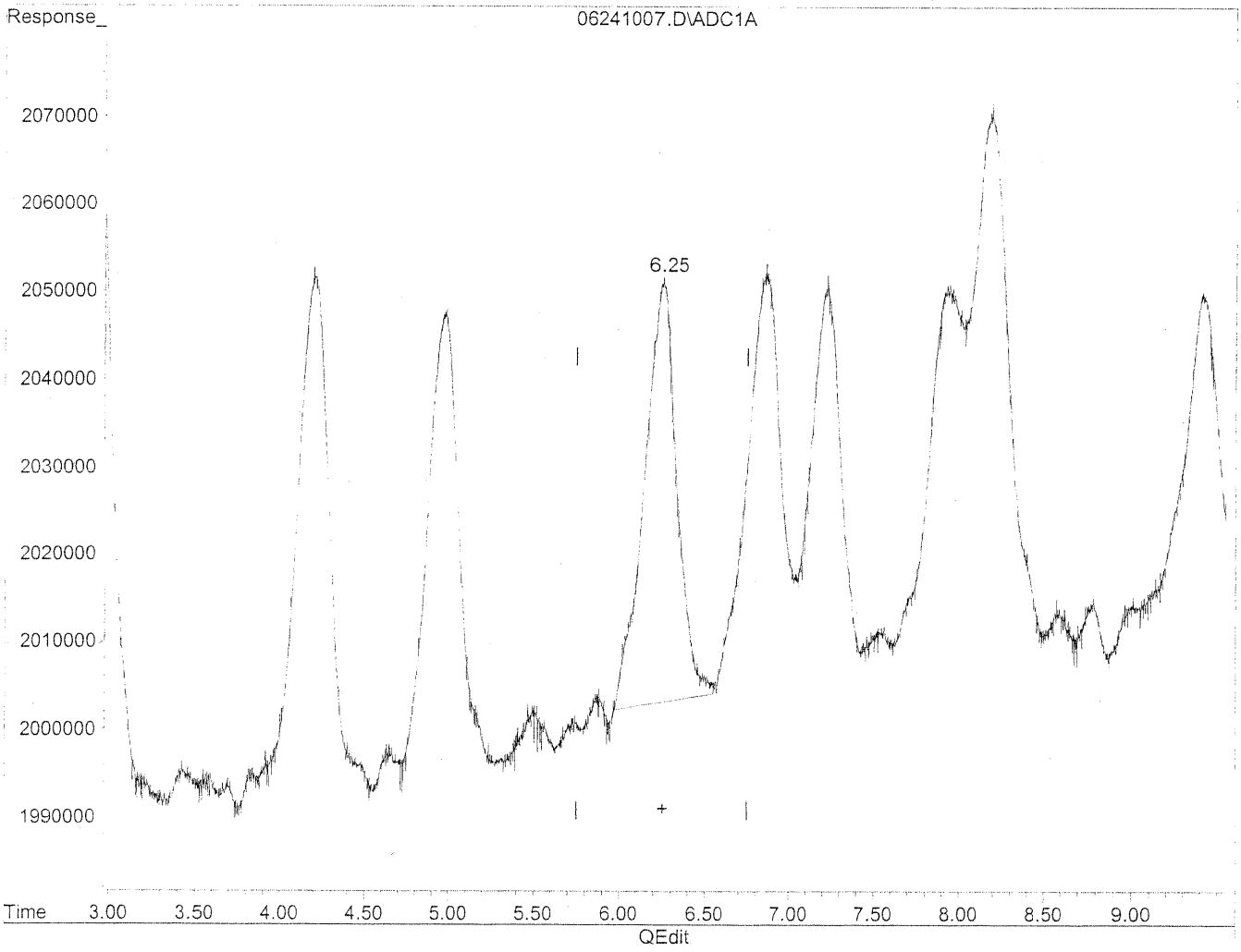
response 7771997

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:39 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.25min 86.389ng/ml m

response 6214018

6/24/10

6/25/10

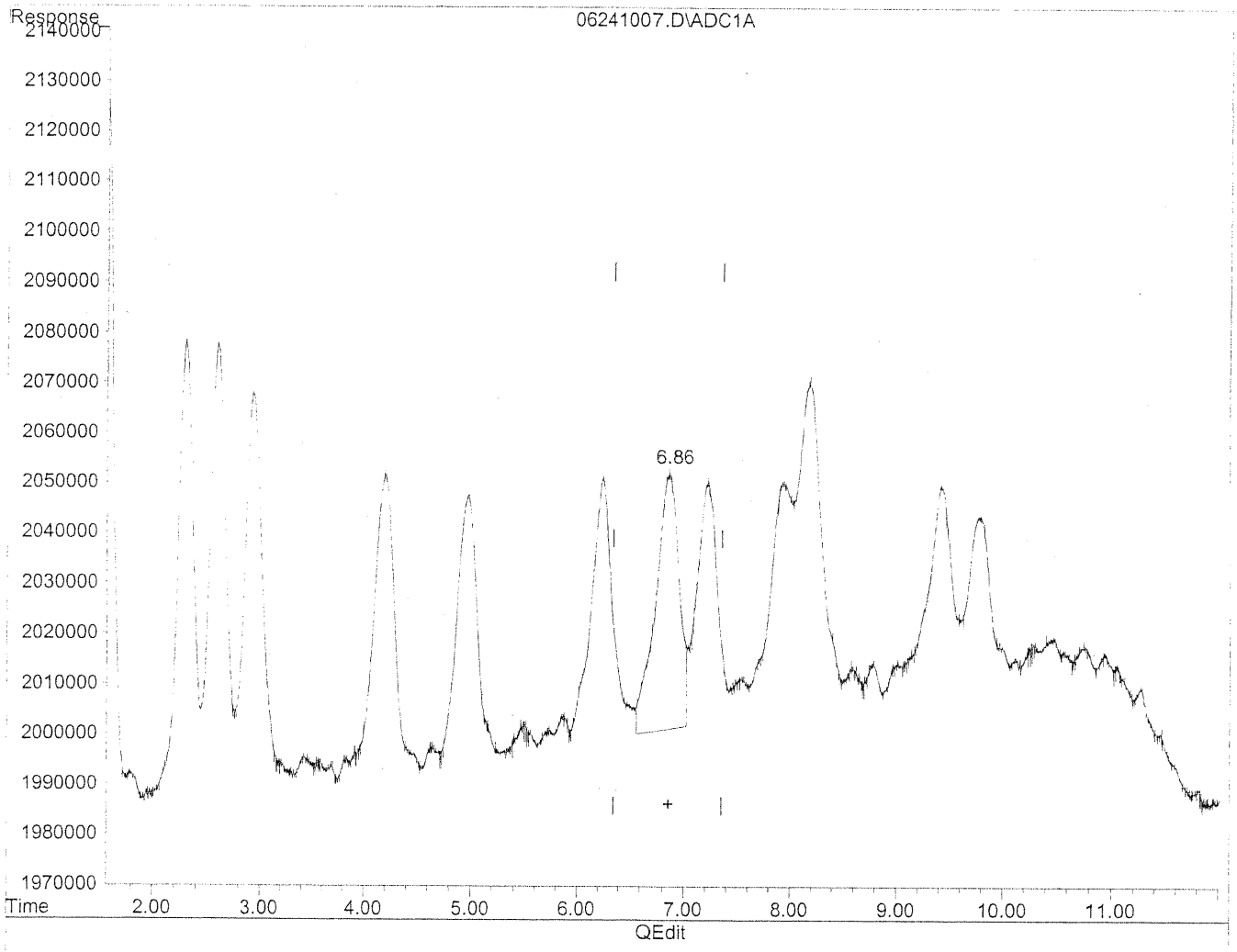
131

(+) = Expected Retention Time
06241007.D TO110610.M Thu Jun 24 16:28:19 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:36 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

6.86min 90.003ng/ml

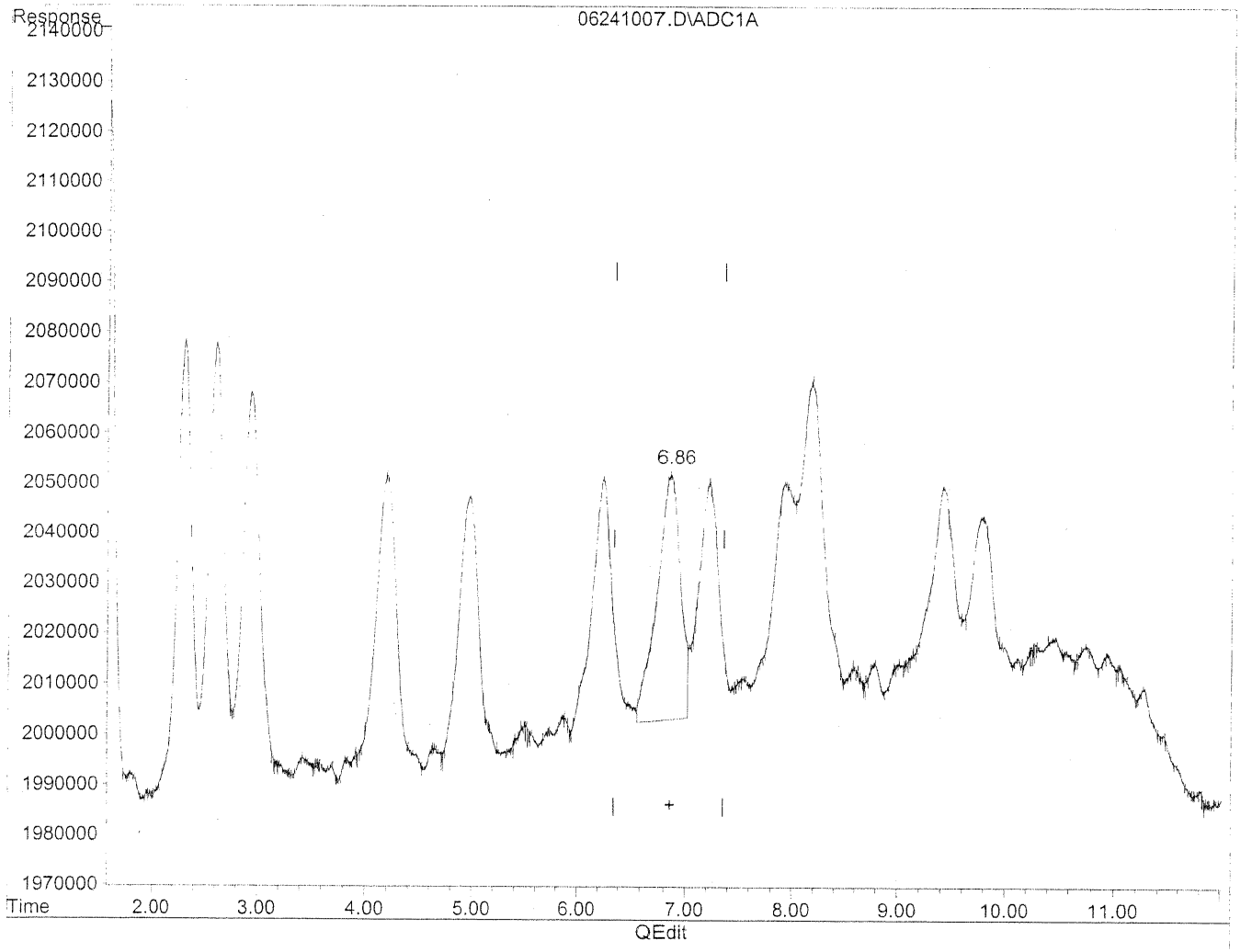
response 7620093

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:36 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

6.86min 82.866ng/ml m

response 7015814

TC 4/28/10

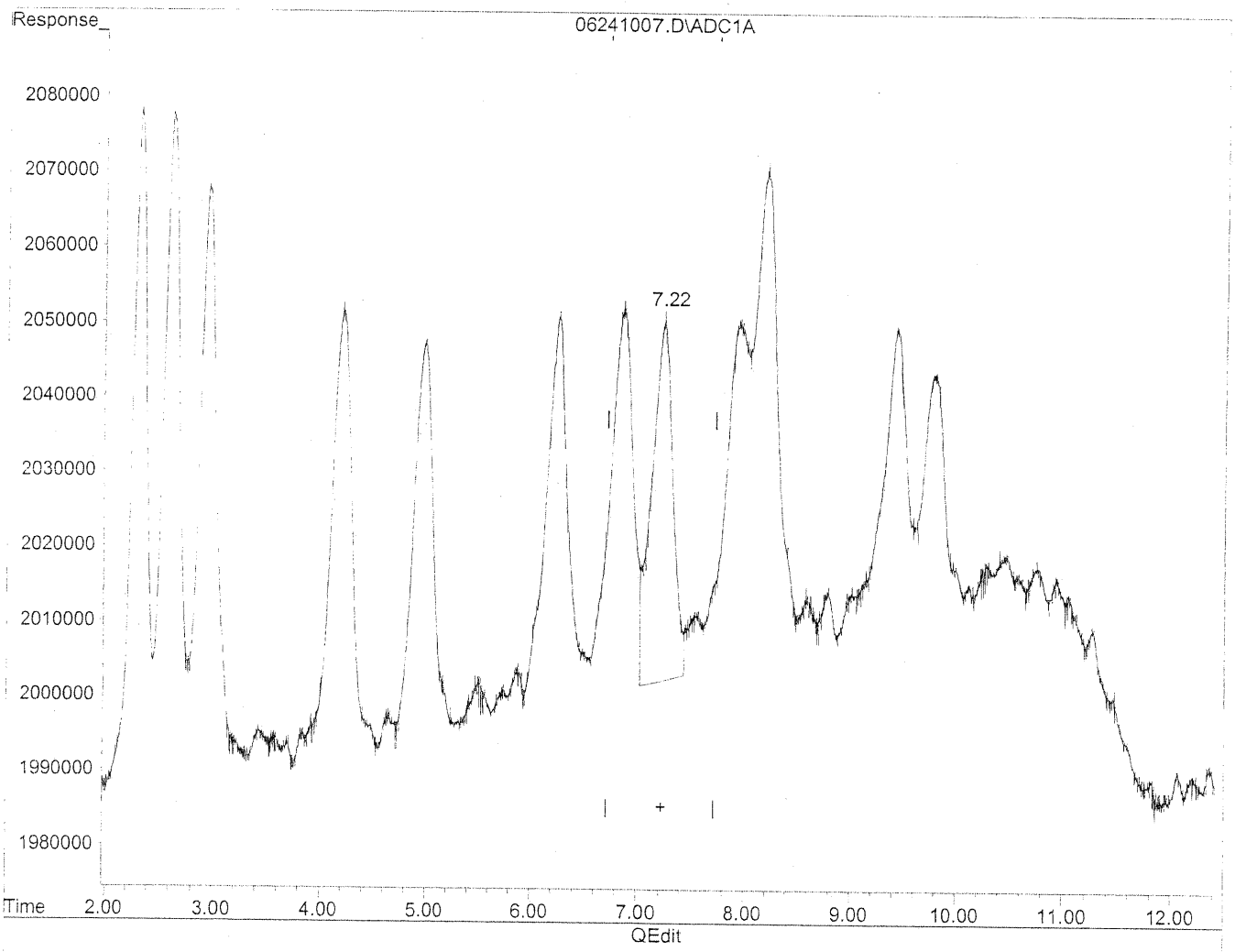
PC 6/25/10

(+) = Expected Retention Time
06241007.D TO110610.M Thu Jun 24 13:37:41 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:36 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.22min 81.357ng/ml

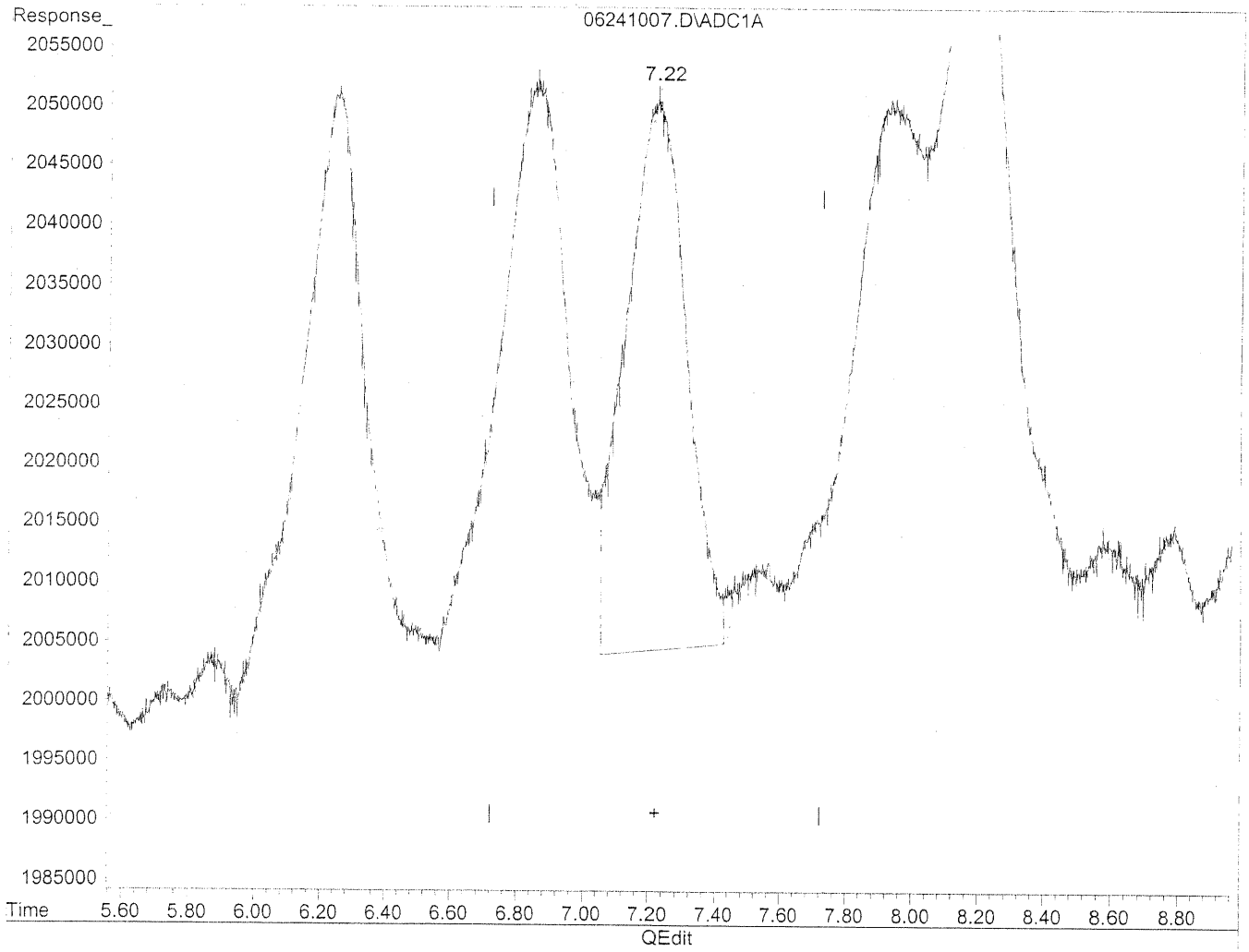
response 6511790

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:39 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.22min 72.167ng/ml m

response 5776234

HL 6/24/10

BL MP 6/25/10

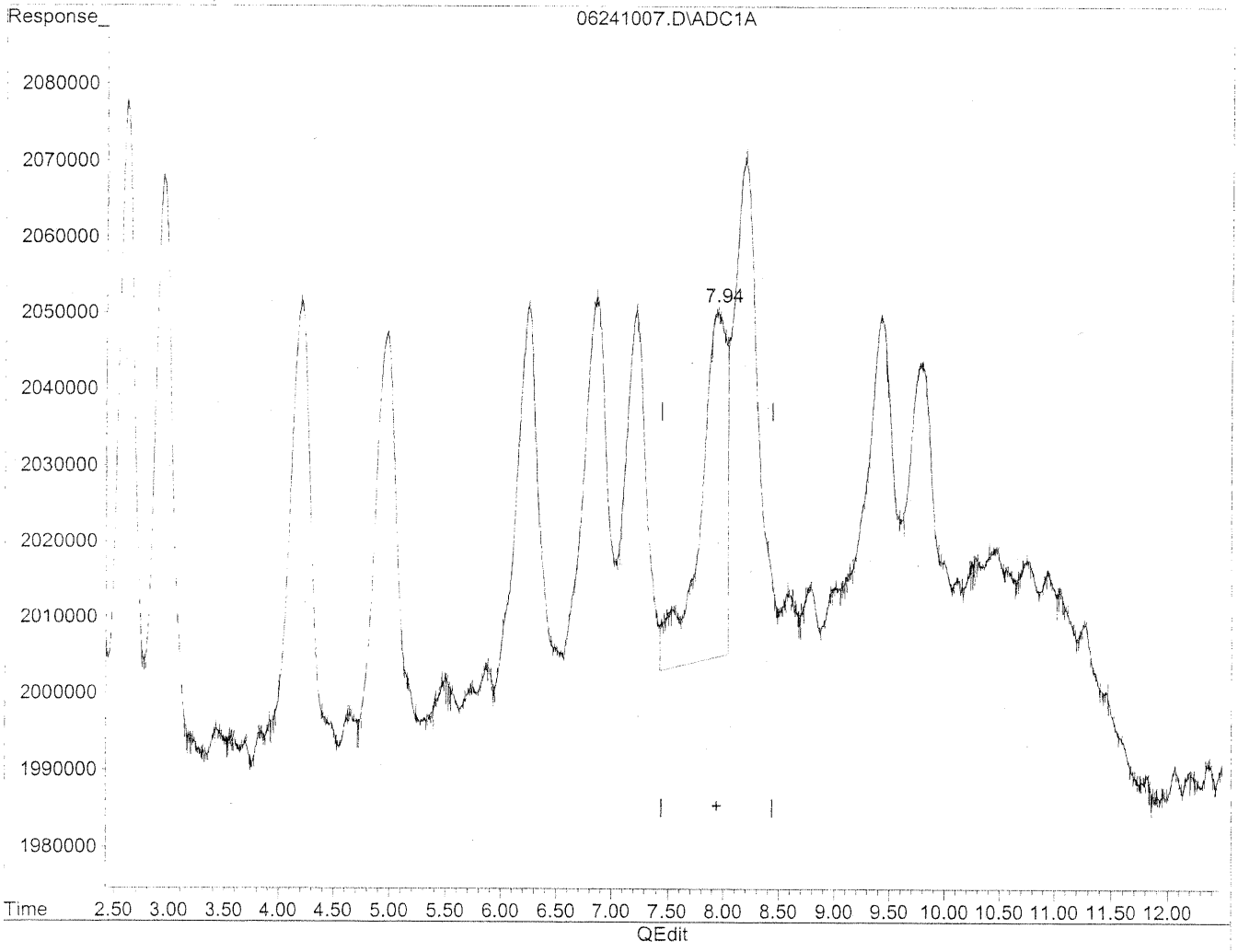
135

(+) = Expected Retention Time
06241007.D TO110610.M Thu Jun 24 16:29:15 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:36 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

7.93min 115.995ng/ml

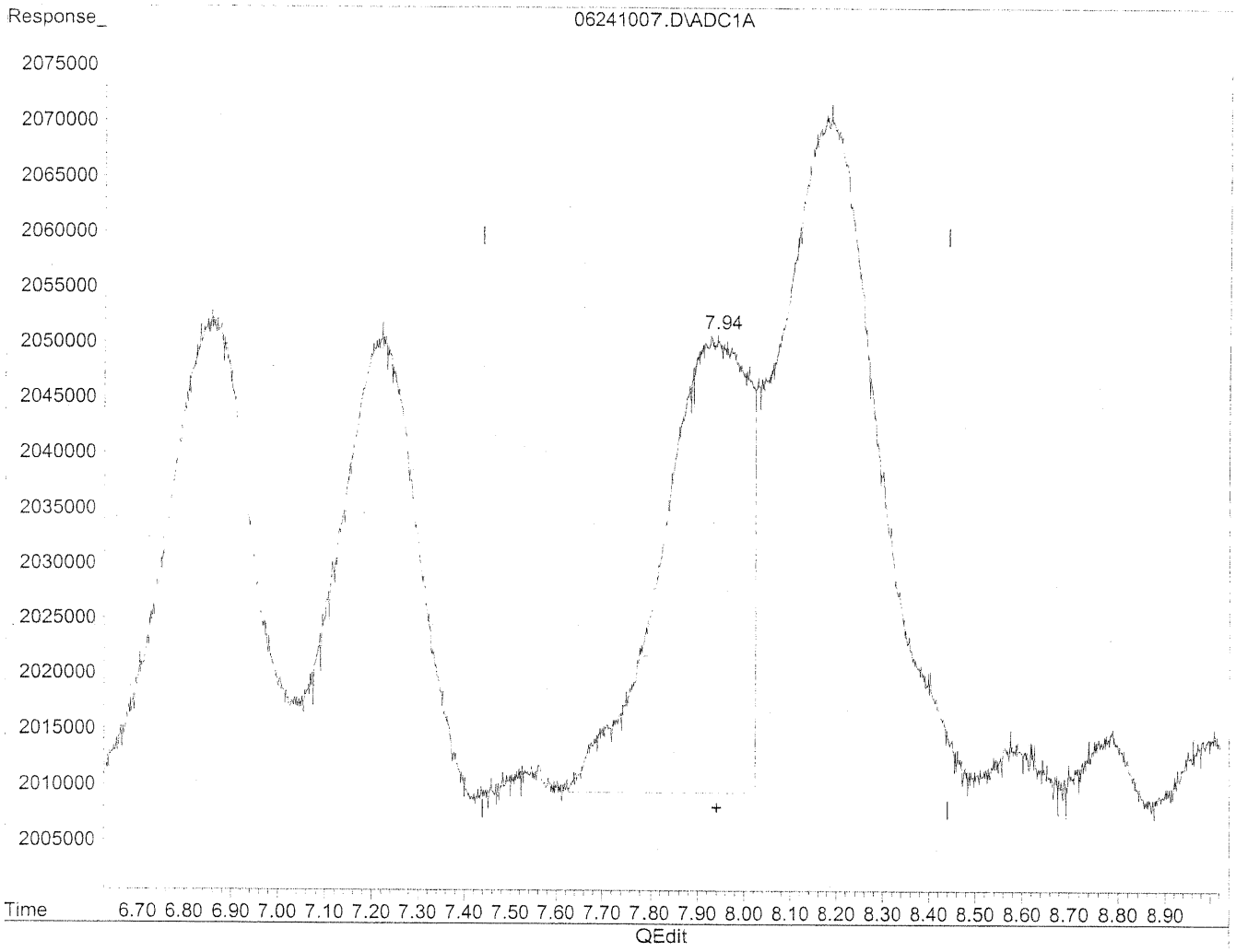
response 7599288

(+) = Expected Retention Time
06241007.D TO110610.M Thu Jun 24 13:38:23 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:49 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 16:48:14 2010
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

7.94min 81.051ng/ml m

response 5309971

HC 6/28/10

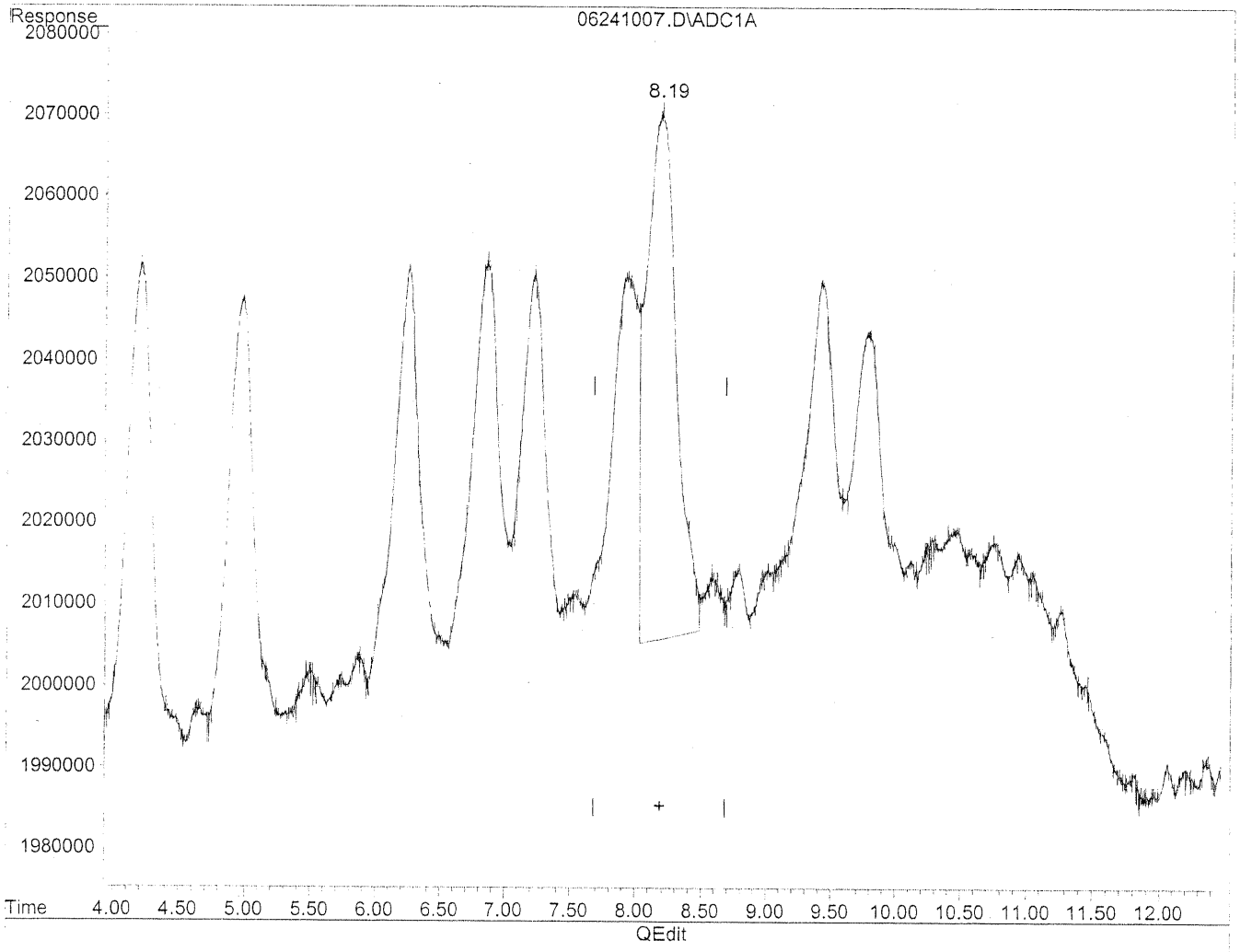
BC MD 6/25/10

(+) = Expected Retention Time
06241007.D TO110610.M Thu Jun 24 16:49:35 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:36 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(10) m,p-Toluaidehyde

8.18min 161.754ng/ml

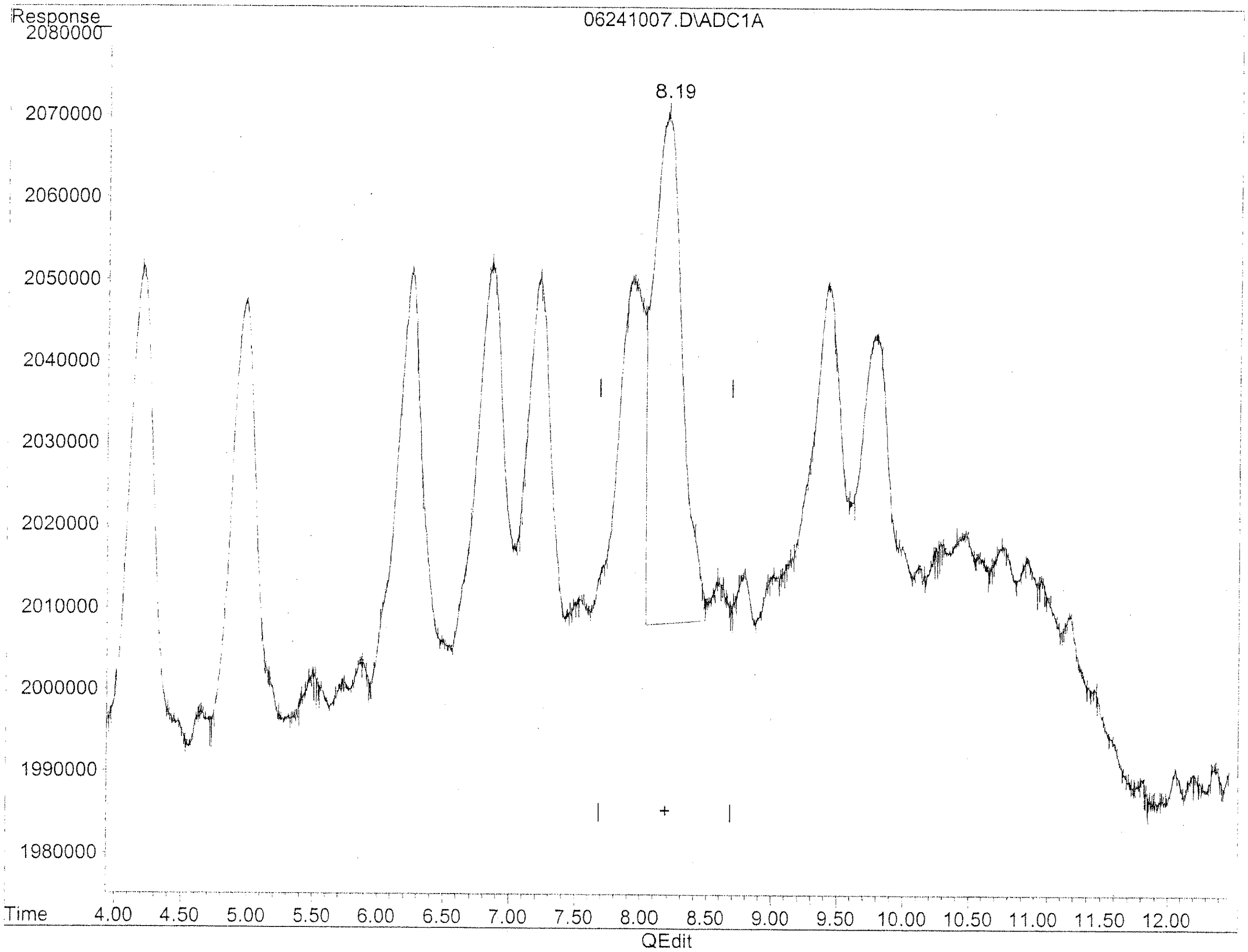
response 9735332

(+) = Expected Retention Time

Quantitation Report

```
Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D      Vial: 5
Acq On    : 24 Jun 2010  12:45                        Operator: MD
Sample    : 100ng/ml TO-11A S21-03091009              Inst    : LC 01
Misc      :                                           Multiplr: 1.00
IntFile   : autoint1.e
Quant Time: Jun 24 13:36 19110  Quant Results File: TO110610.RES

Method          : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title           : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update     : Thu Jun 24 13:26:29 2010
Response via    : Multiple Level Calibration
```



(10) m,p-Tolualdehyde

8.19min 149.541ng/ml m

response 9000241

JK 6/28/10

BV MD 6/25/10

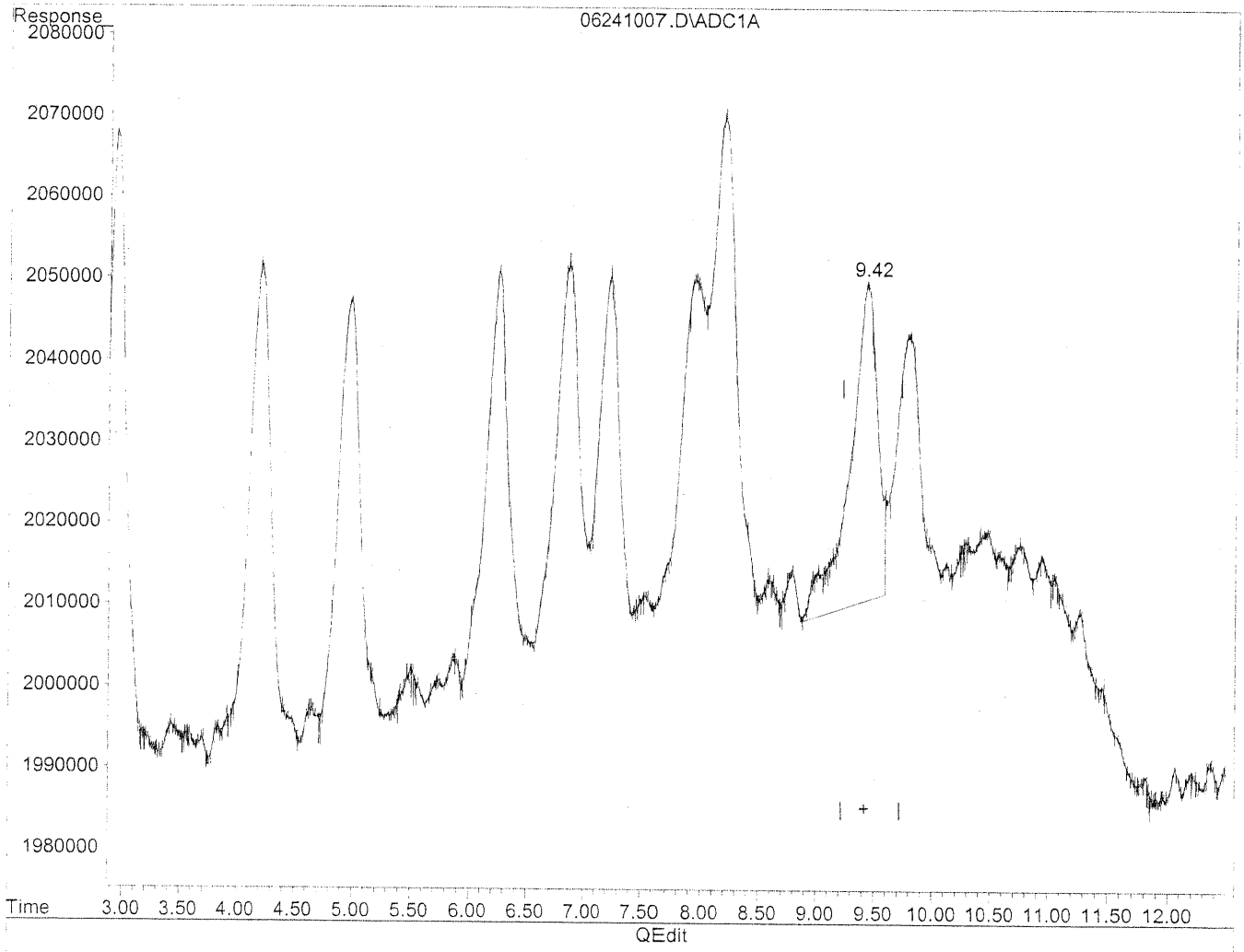
(+) = Expected Retention Time

06241007.D   TO110610.M   Thu Jun 24 13:38:46 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:36 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(11) Hexaldehyde

9.42min 83.691ng/ml

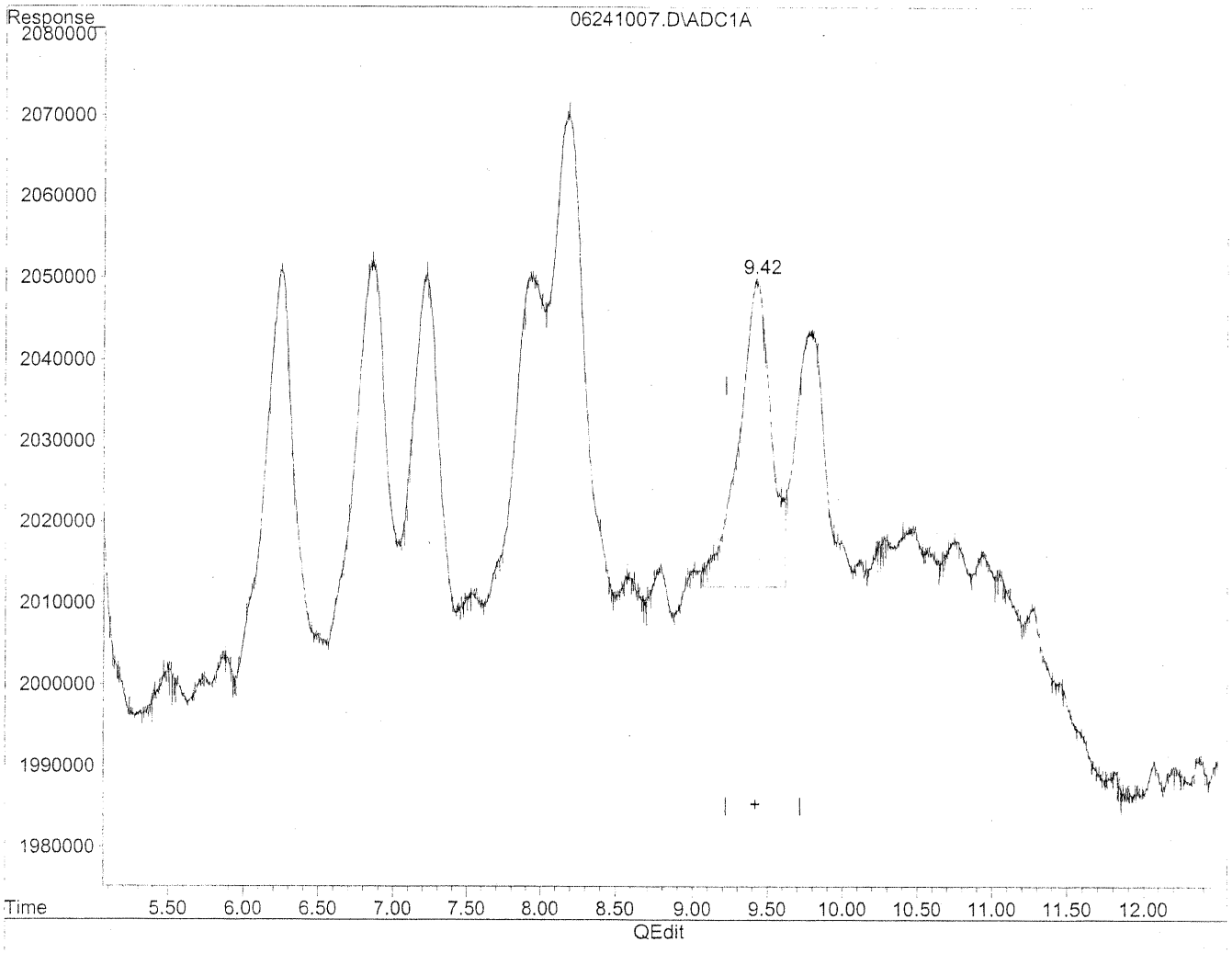
response 6368954

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:49 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
Response via : Multiple Level Calibration



(11) Hexaldehyde

9.42min 73.979ng/ml m

response 5629872

6/28/10

6/28/10

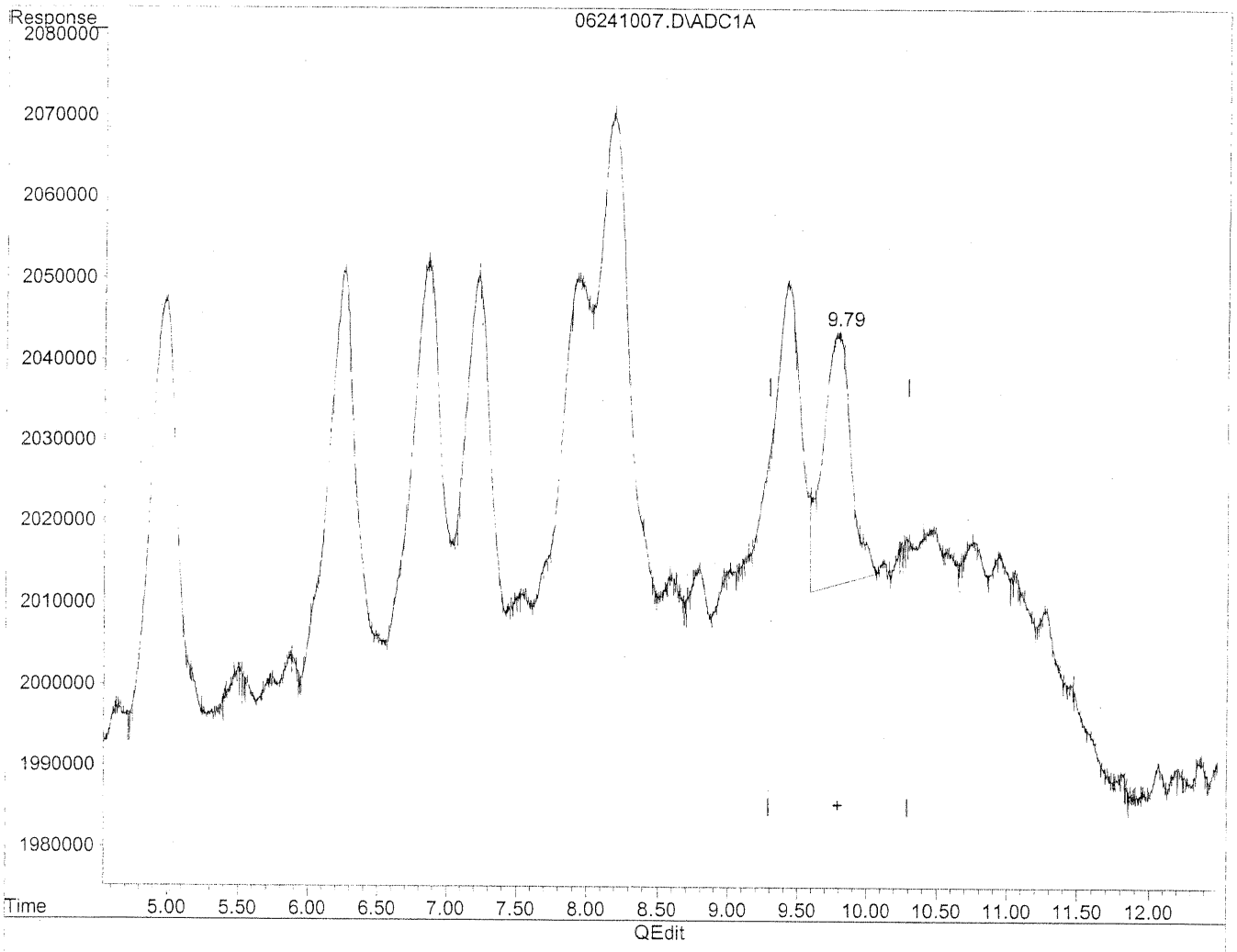
141

(+) = Expected Retention Time
06241007.D TO110610.M Fri Jun 25 10:21:48 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:36 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

9.78min 84.721ng/ml

response 4442936

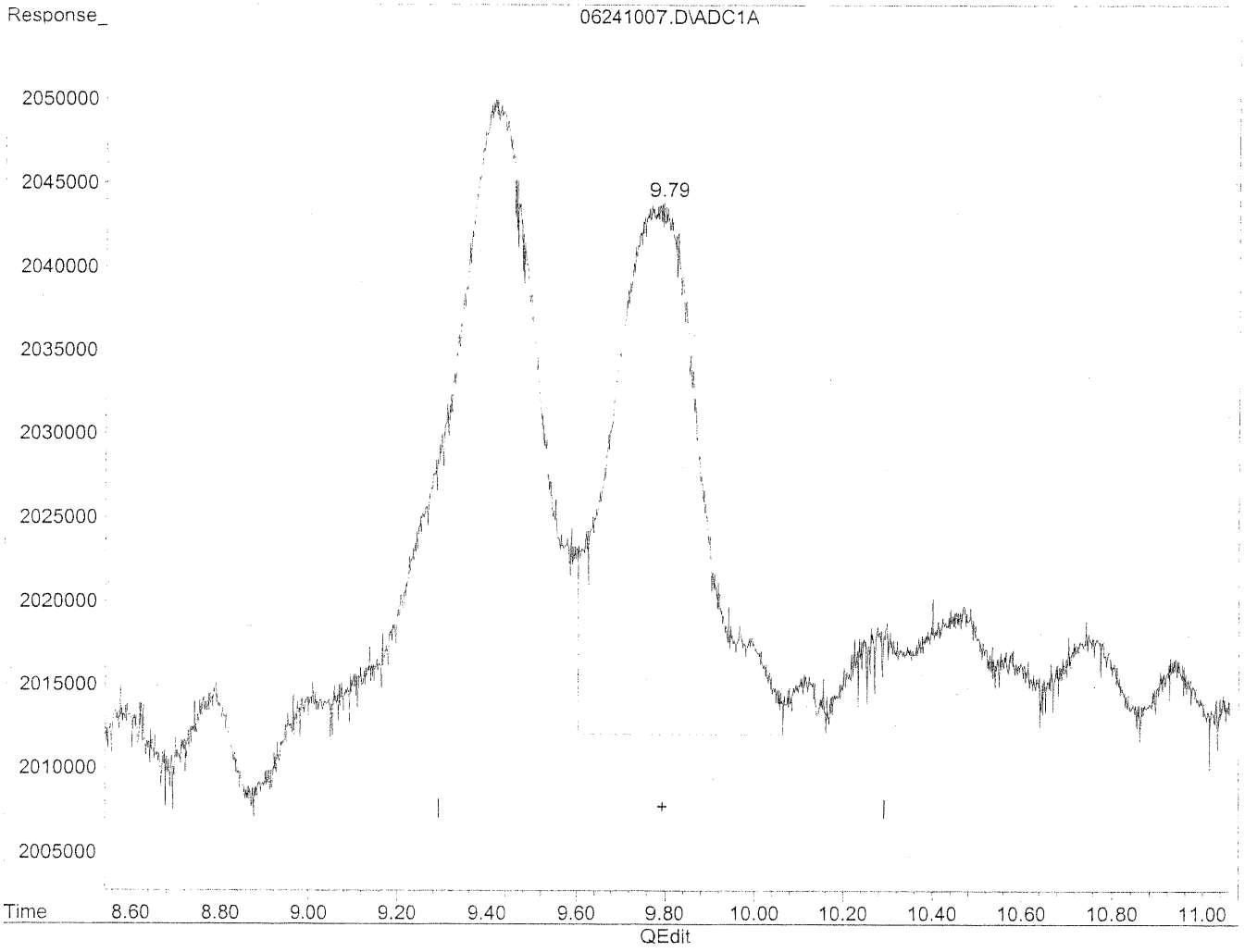
142

(+) = Expected Retention Time
06241007.D TO110610.M Thu Jun 24 13:39:18 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241007.D Vial: 5
Acq On : 24 Jun 2010 12:45 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:49 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 16:48:14 2010
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

9.79min 85.755ng/ml m

response 4497151

HL 6/24/10

TC 6/25/10

143

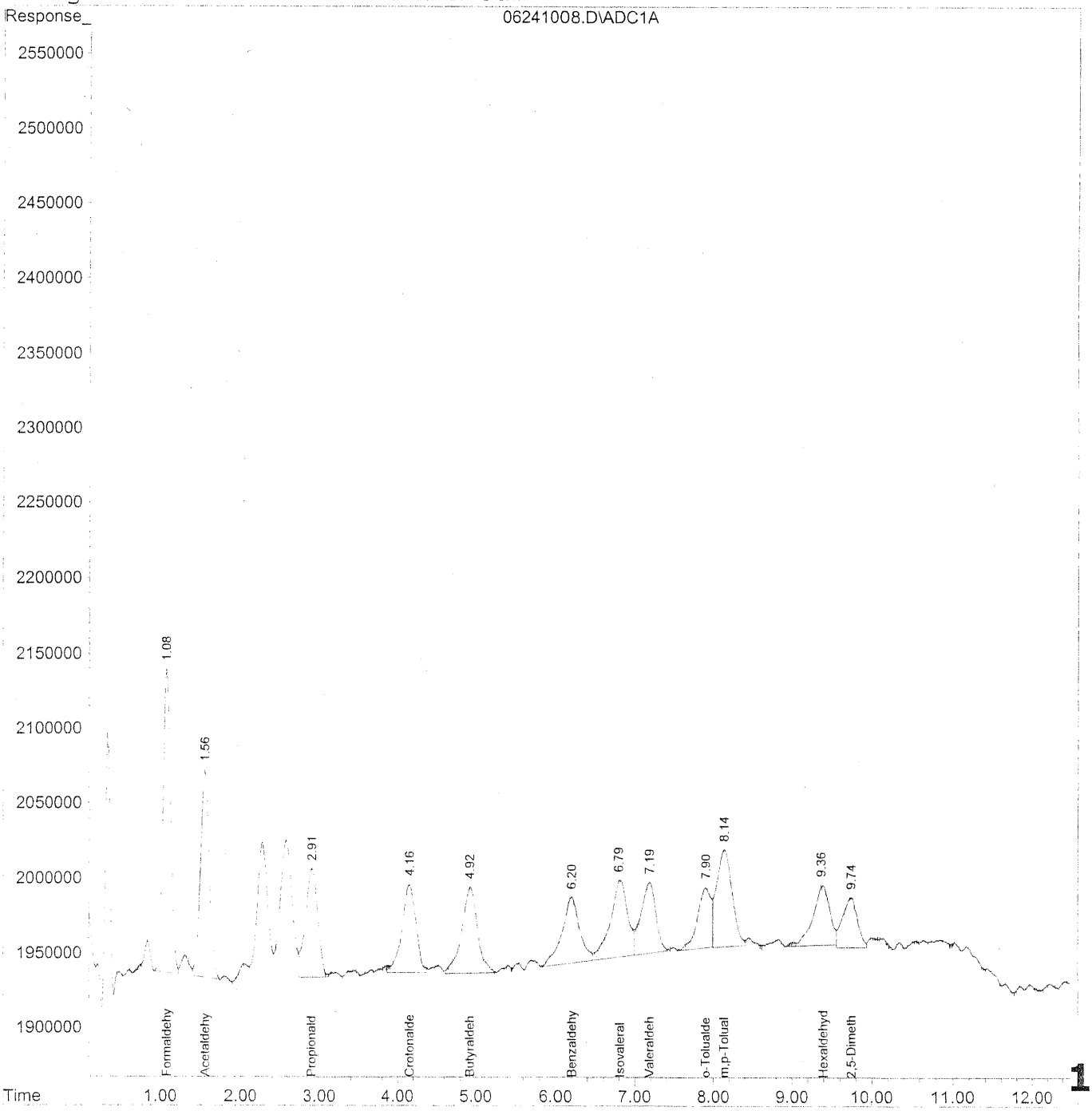
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241008.D Vial: 6
Acq On : 24 Jun 2010 12:58 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:50 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
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\2010_06\24\06241008.D Vial: 6
Acq On : 24 Jun 2010 12:58 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:50 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
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

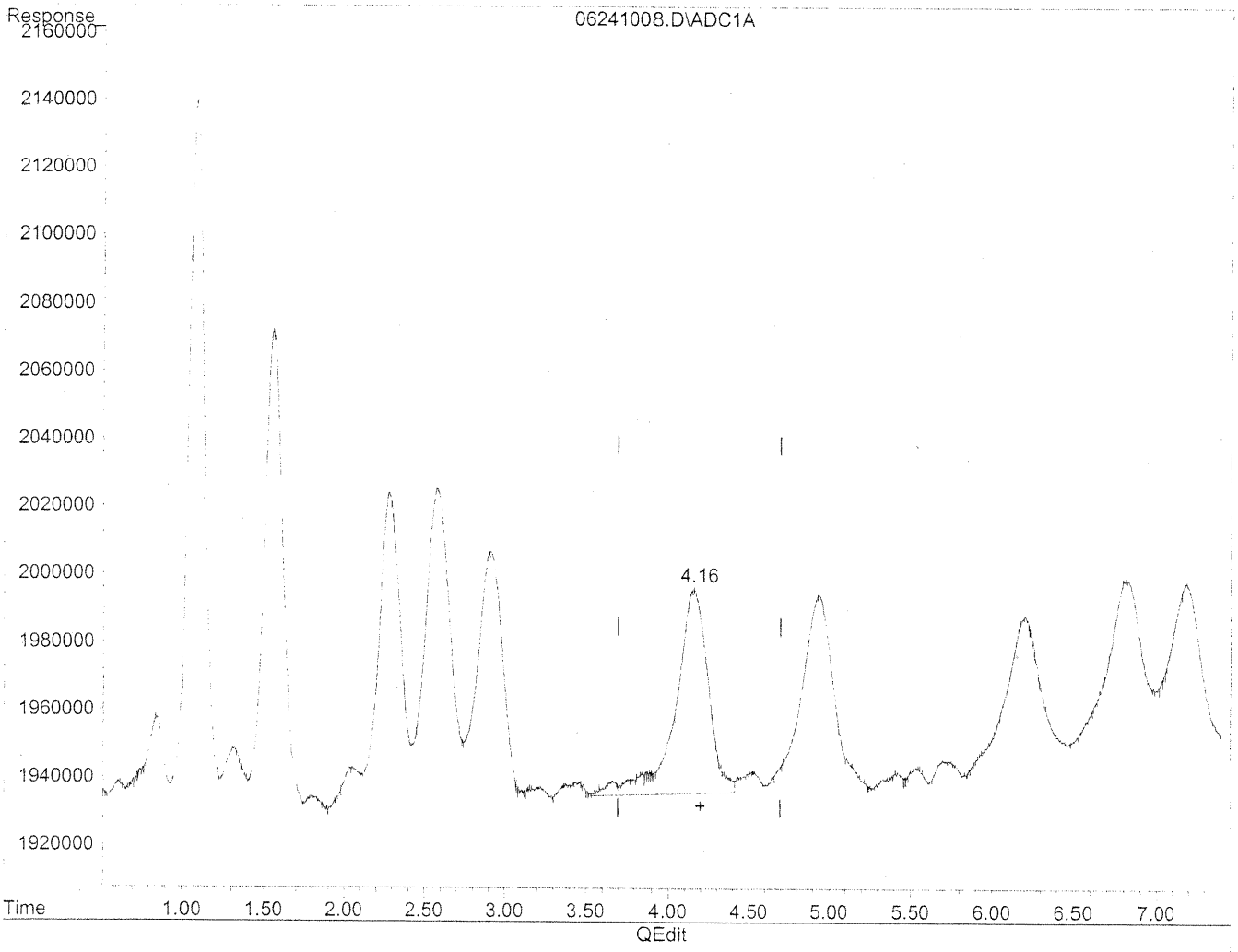
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.08	13511387	71.138	ng/ml
2)	Acetaldehyde	1.55	10740786	73.557	ng/ml
3)	Propionaldehyde	2.91	8027964	72.404	ng/ml
4)	Crotonaldehyde	4.16	7500614	75.147	ng/mlm
5)	Butyraldehyde	4.93	8049959	87.157	ng/ml
6)	Benzaldehyde	6.20	6787382	97.676	ng/ml
7)	Isovaleraldehyde	6.81	7873759	97.612	ng/ml
8)	Valeraldehyde	7.19	6332528	84.989	ng/ml
9)	o-Tolualdehyde	7.90	4800335	74.257	ng/mlm
10)	m,p-Tolualdehyde	8.14	8786398	154.715	ng/mlm
11)	Hexaldehyde	9.39	5891760	82.955	ng/ml
12)	2,5-Dimethylbenzaldehyde	9.74	4437325	89.918	ng/mlm

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241008.D Vial: 6
Acq On : 24 Jun 2010 12:58 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:40 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.15min 86.940ng/ml

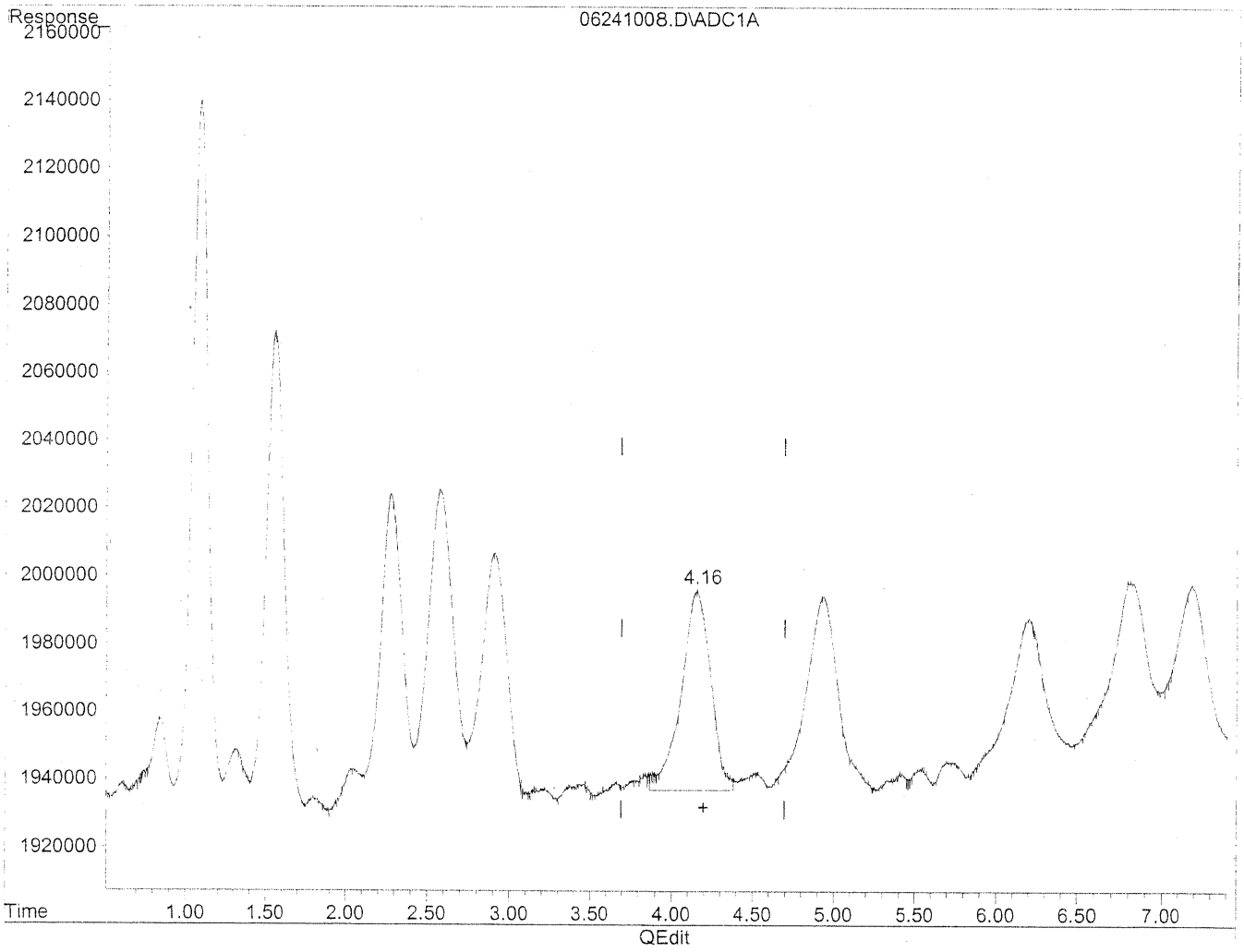
response 8677730

(+) = Expected Retention Time
06241008.D TO110610.M Thu Jun 24 13:40:12 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241008.D Vial: 6
Acq On : 24 Jun 2010 12:58 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:40 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.16min 75.147ng/ml m

response 7500614

HC 6/25/10

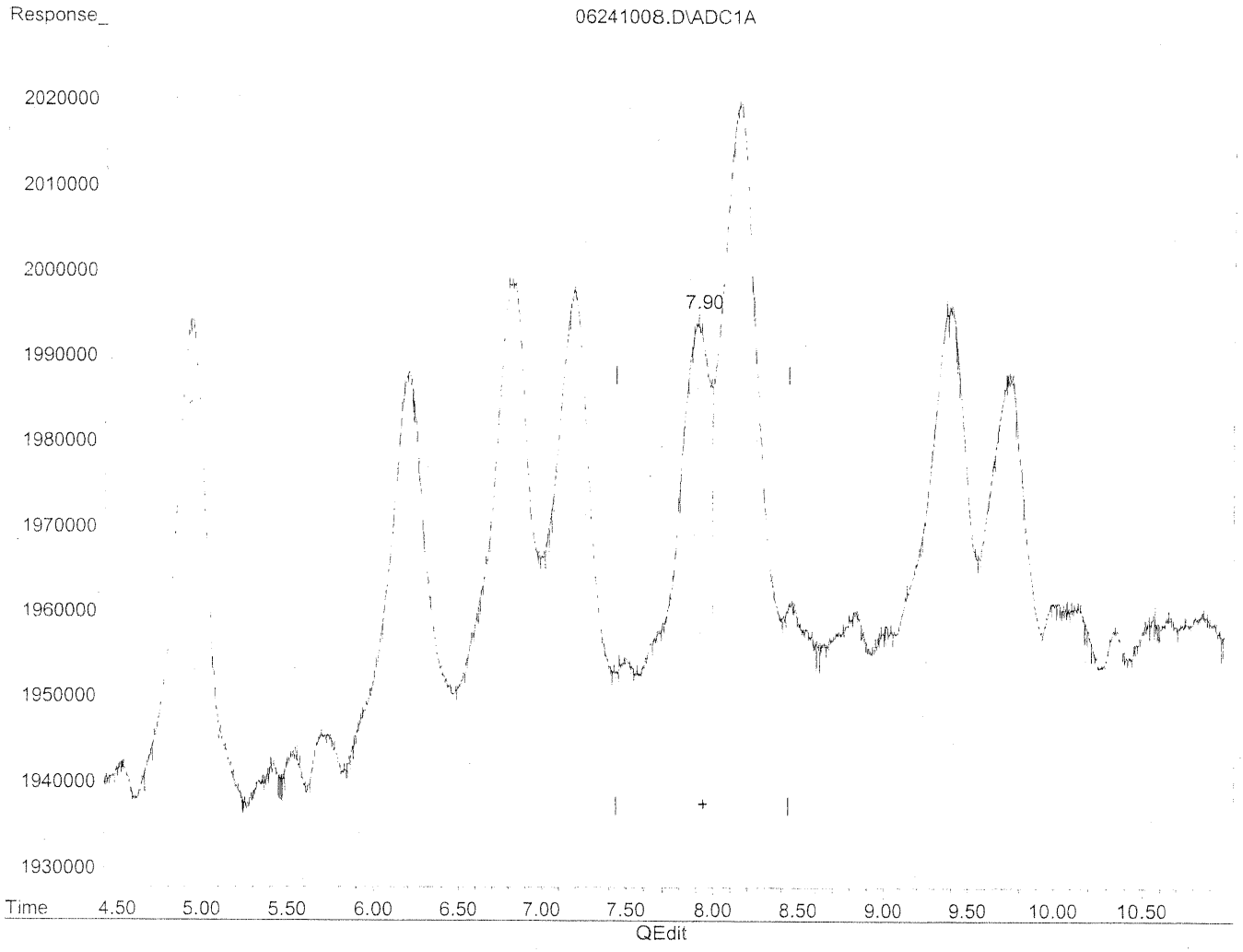
TC 6/25/10

(+) = Expected Retention Time
06241008.D TO110610.M Thu Jun 24 13:40:17 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241008.D Vial: 6
Acq On : 24 Jun 2010 12:58 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:40 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

7.90min 70.602ng/ml

response 4564084

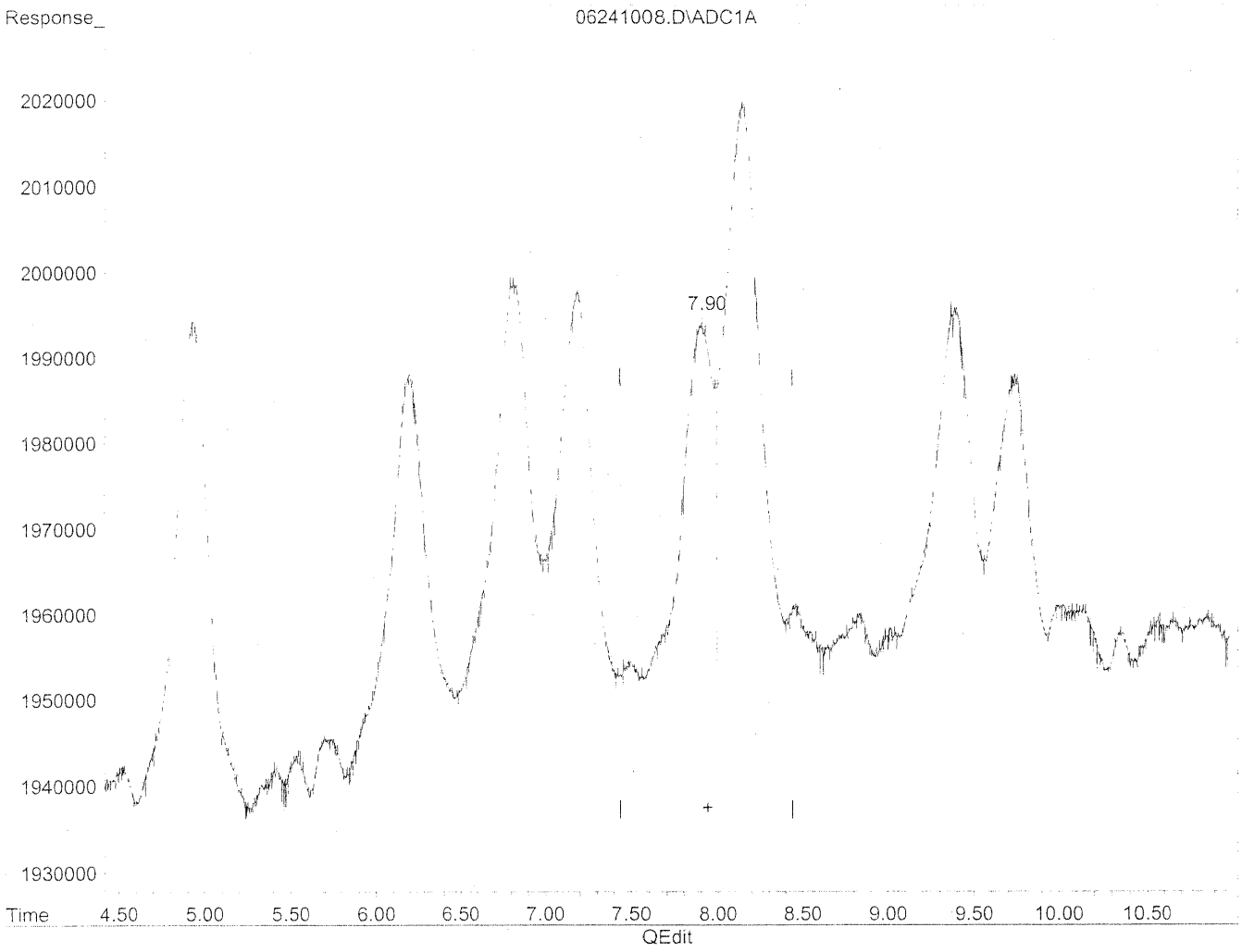
148

(+) = Expected Retention Time
06241008.D TO110610.M Thu Jun 24 13:40:42 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241008.D Vial: 6
Acq On : 24 Jun 2010 12:58 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:40 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

7.90min 74.257ng/ml m

response 4800335

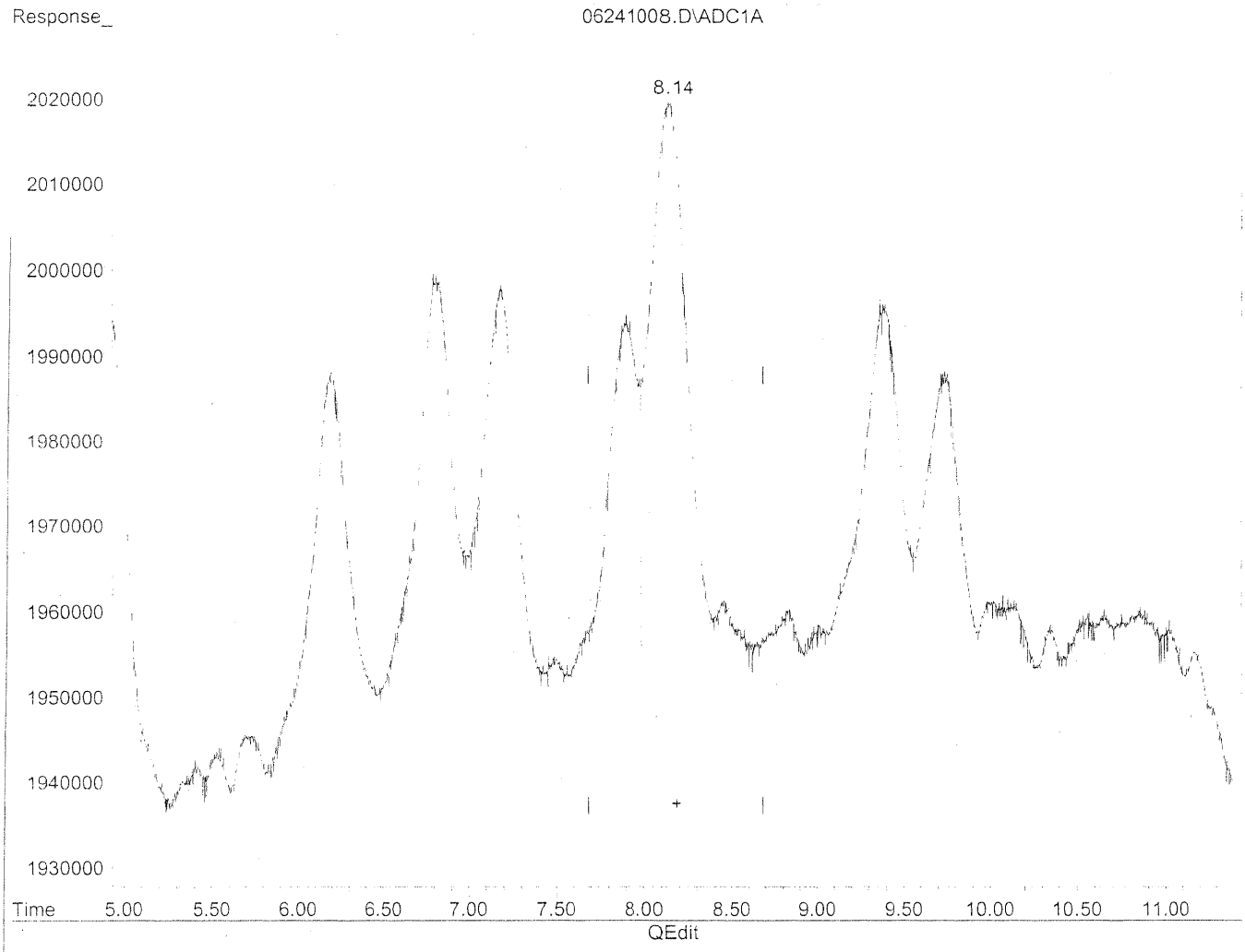
149

(+) = Expected Retention Time
06241008.D TO110610.M Thu Jun 24 13:40:52 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241008.D Vial: 6
Acq On : 24 Jun 2010 12:58 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:40 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

8.14min 147.192ng/ml

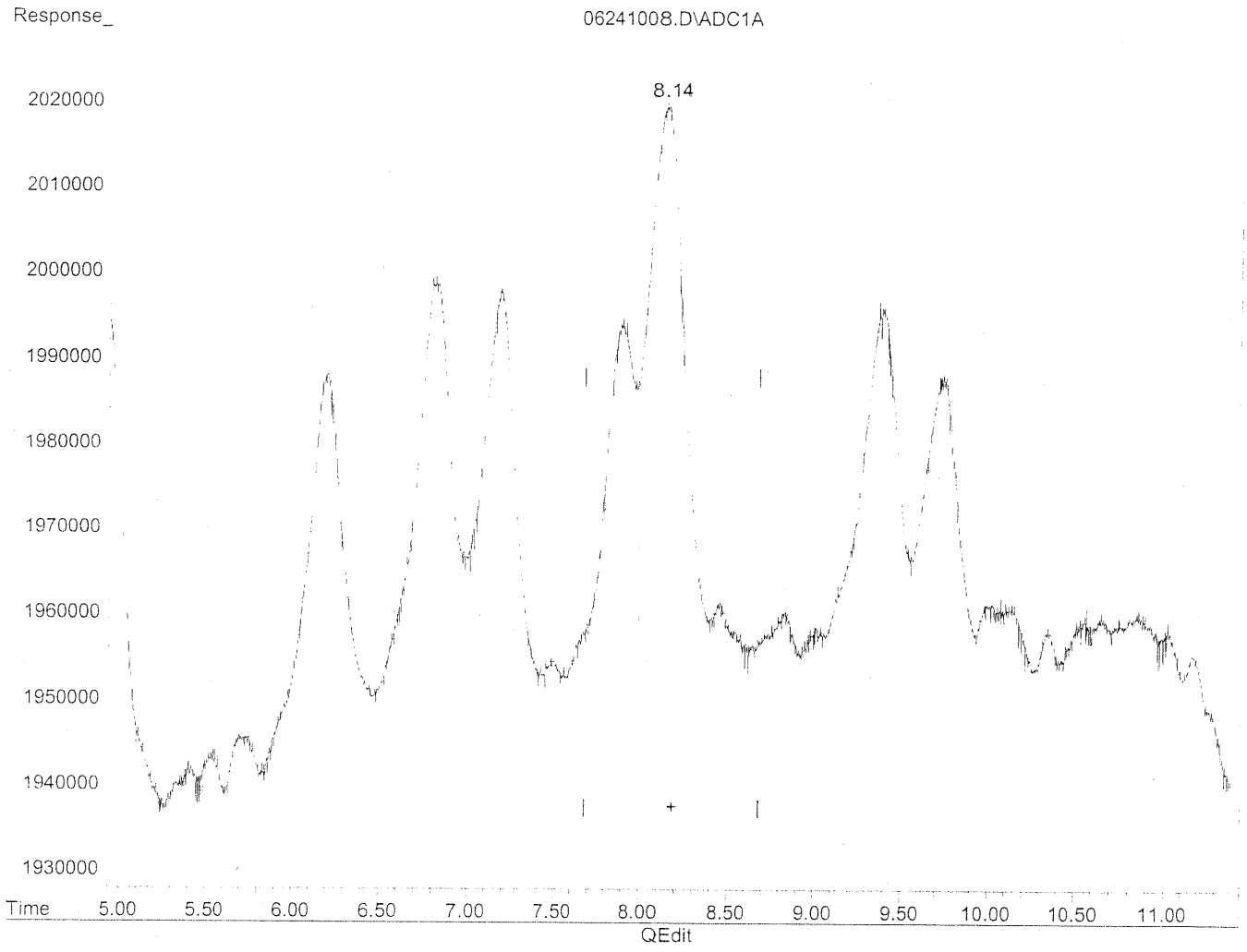
response 8359189

(+) = Expected Retention Time
06241008.D TO110610.M Thu Jun 24 13:40:58 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241008.D Vial: 6
Acq On : 24 Jun 2010 12:58 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:40 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

8.14min 154.715ng/ml m

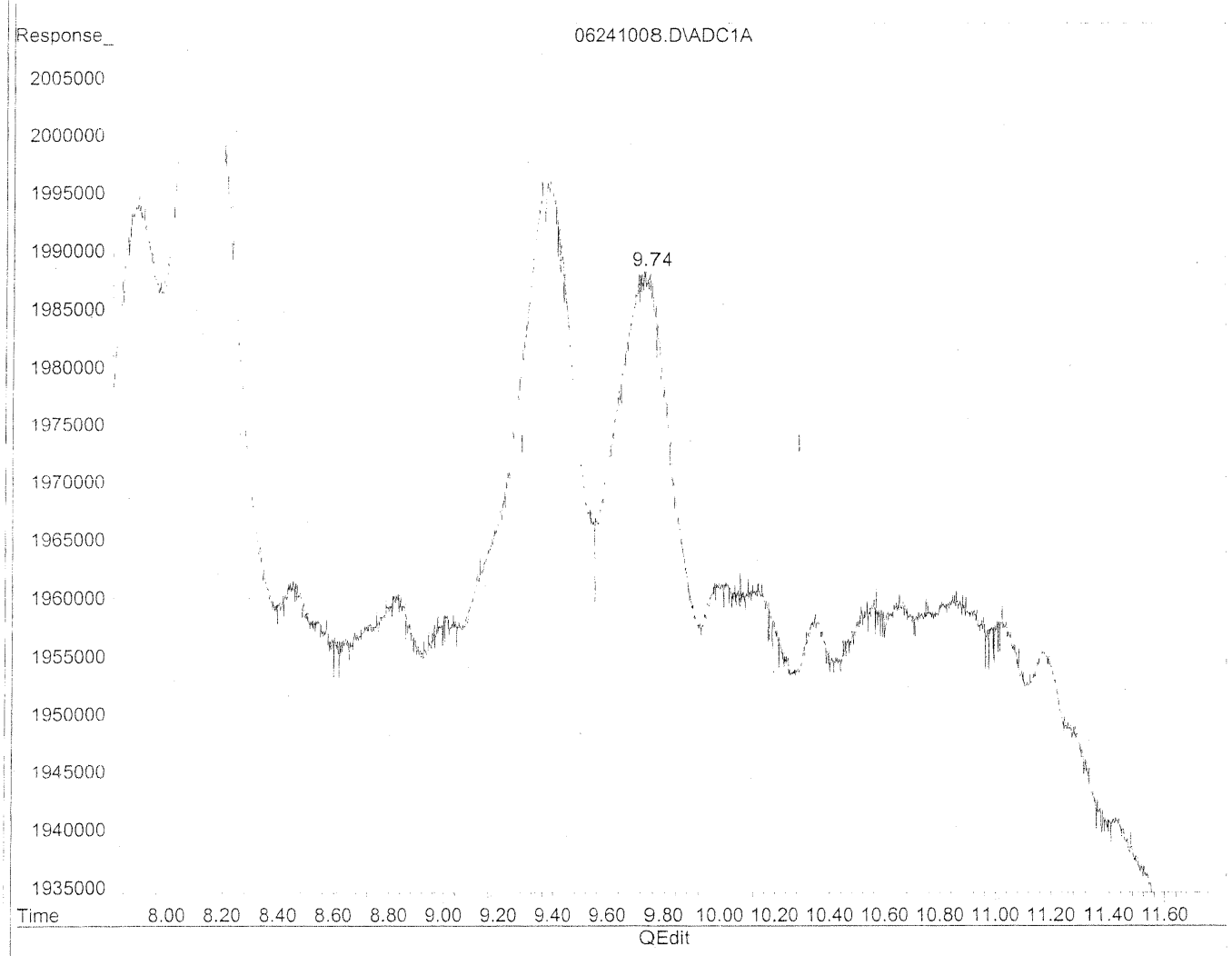
response 8786398

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241008.D Vial: 6
Acq On : 24 Jun 2010 12:58 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:41 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 16:30:28 2010
Response via : Multiple Level Calibration



(12) 2.5-Dimethylbenzaldehyde

9.74min 78.272ng/ml

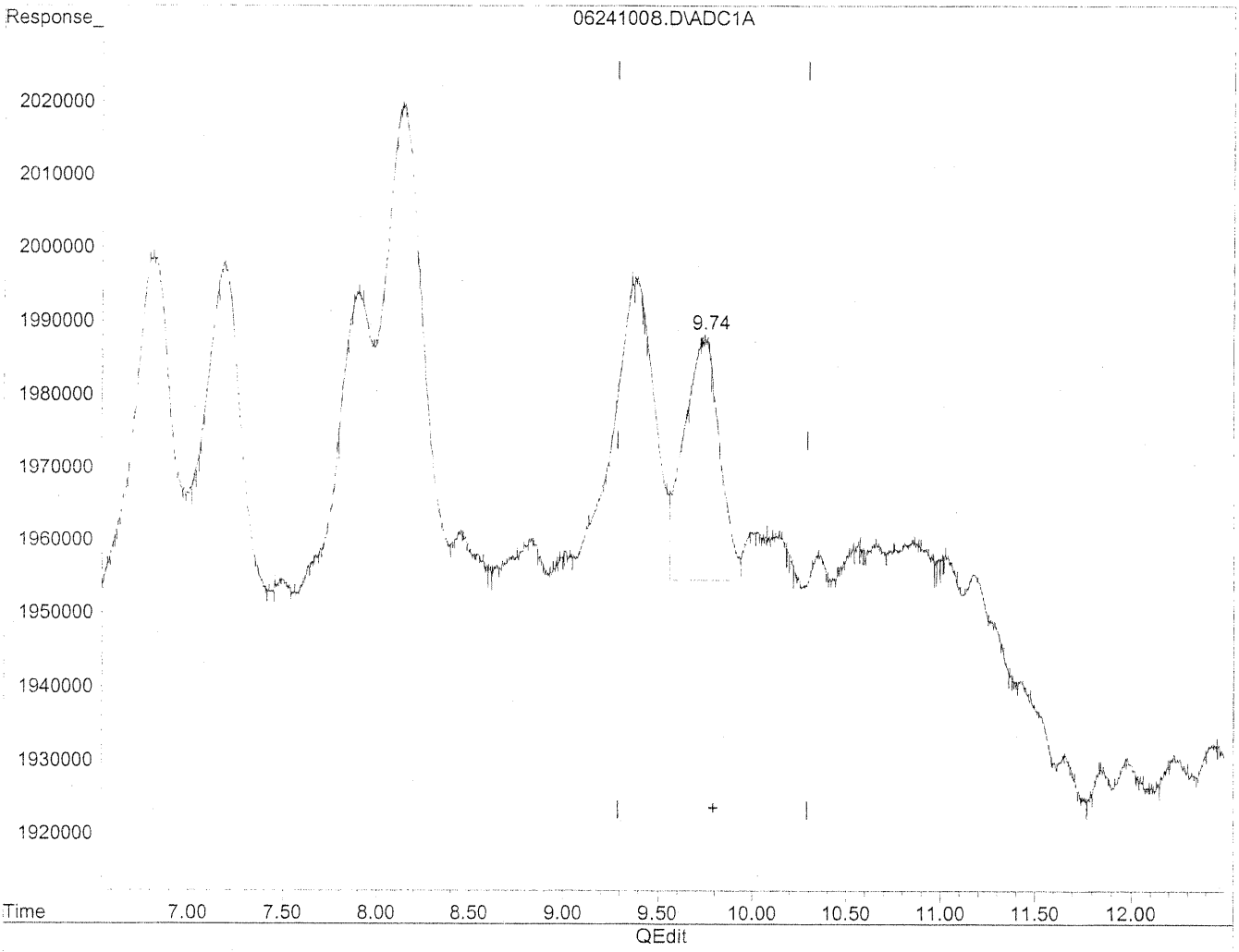
response 3862623

(+) = Expected Retention Time
06241008.D TO110610.M Thu Jun 24 16:31:11 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241008.D Vial: 6
Acq On : 24 Jun 2010 12:58 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:31 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 16:49:50 2010
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

9.74min 89.918ng/ml m

response 4437325

6/25/10

6/25/10

153

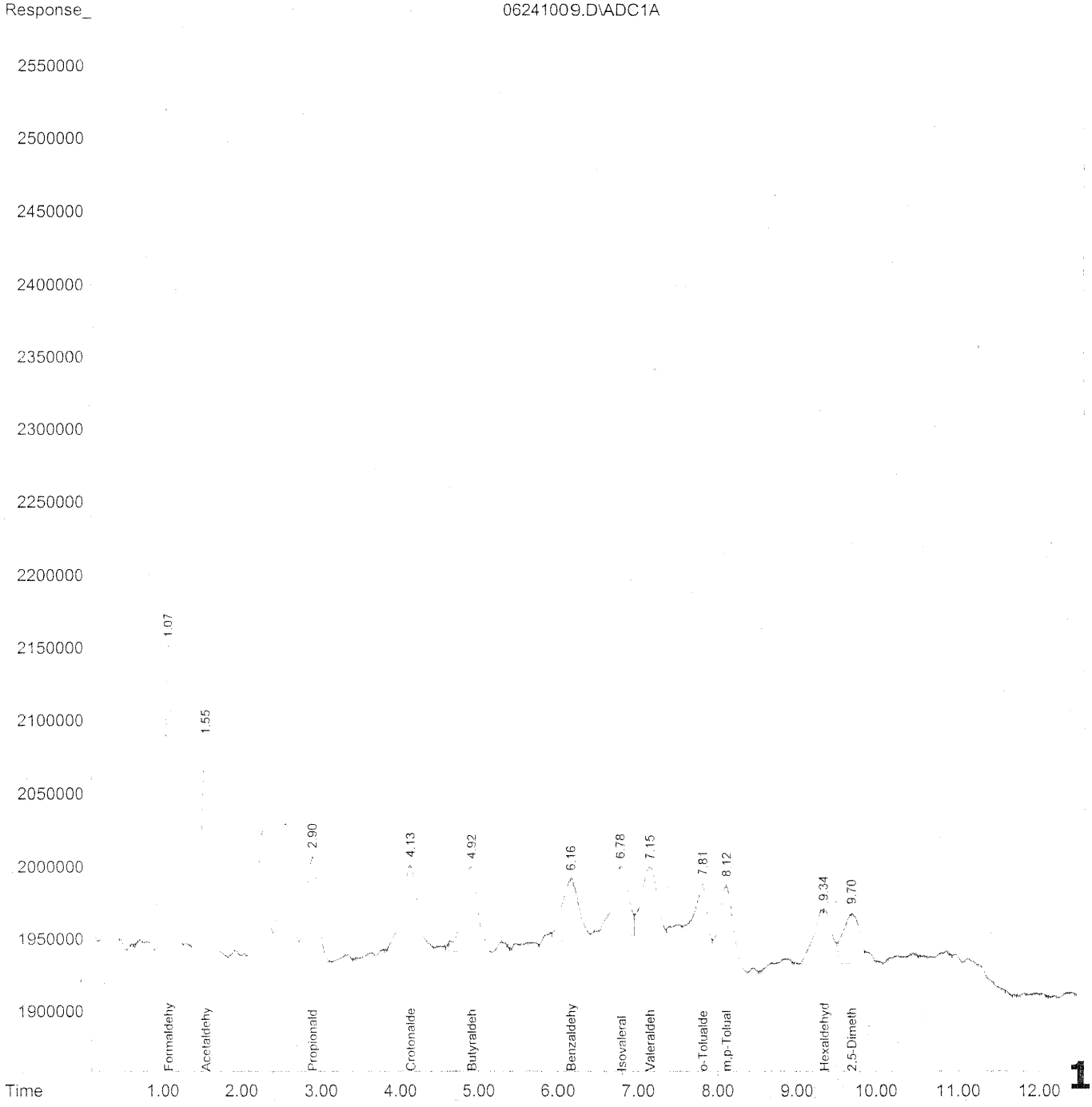
(+) = Expected Retention Time
06241008.D TO110610.M Thu Jun 24 16:50:40 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:47 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
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\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:47 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
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

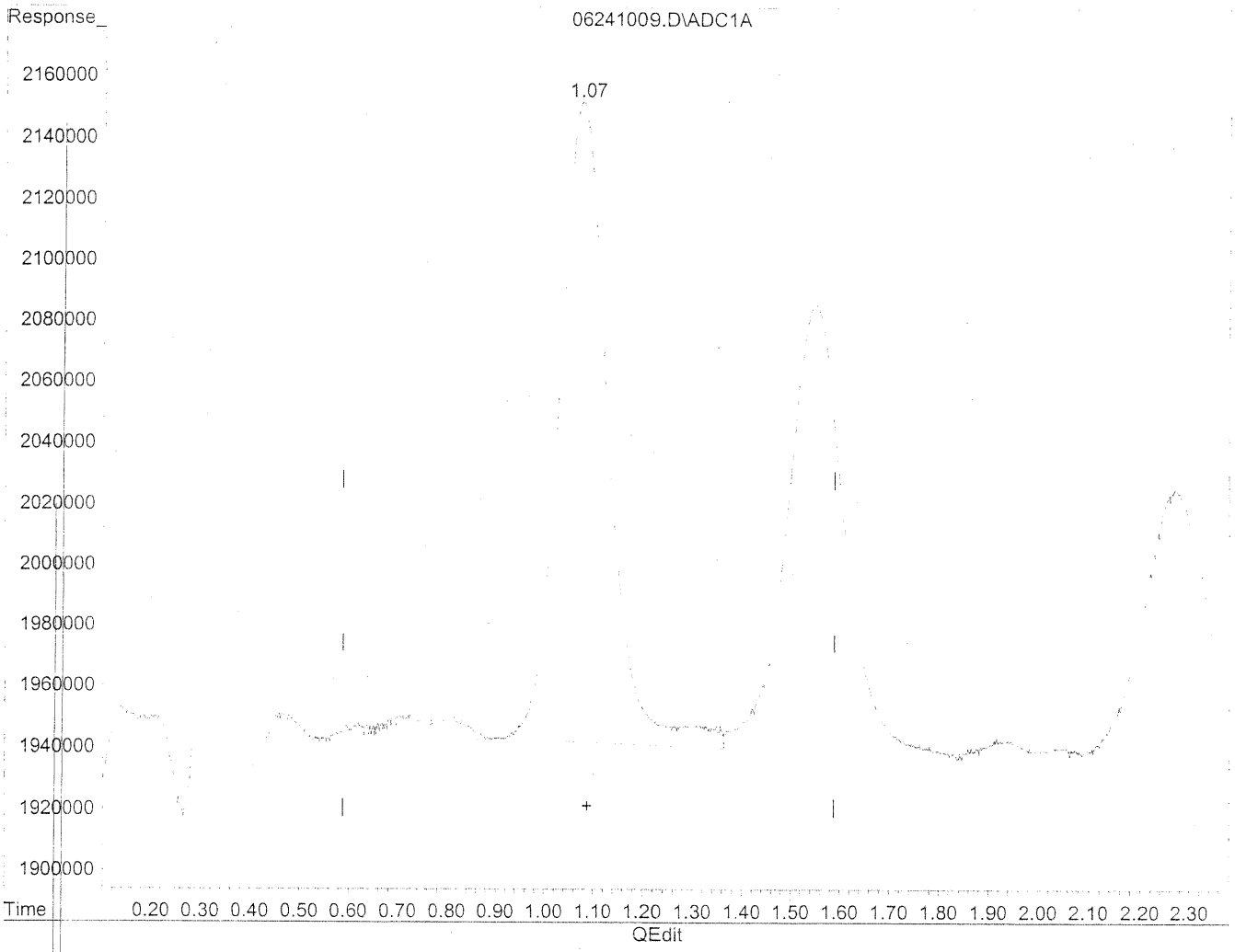
	Compound	R.T.	Response	Conc	Units
Target Compounds					
1)	Formaldehyde	1.07	14186865	74.751	ng/mlm
2)	Acetaldehyde	1.55	11127714	76.129	ng/mlm
3)	Propionaldehyde	2.90	7835935	71.146	ng/ml
4)	Crotonaldehyde	4.13	8490383	85.389	ng/mlm
5)	Butyraldehyde	4.92	8096555	86.753	ng/mlm
6)	Benzaldehyde	6.16	5893577	84.678	ng/mlm
7)	Isovaleraldehyde	6.79	6729977	82.699	ng/mlm
8)	Valeraldehyde	7.15	6646232	88.470	ng/mlm
9)	o-Tolualdehyde	7.81	4566176	71.724	ng/mlm
10)	m,p-Tolualdehyde	8.12	8138944	143.539	ng/mlm
11)	Hexaldehyde	9.34	5878688	82.359	ng/mlm
12)	2,5-Dimethylbenzaldehyde	9.70	4860735	98.946	ng/mlm

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:41 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 77.040ng/ml

response 14621253

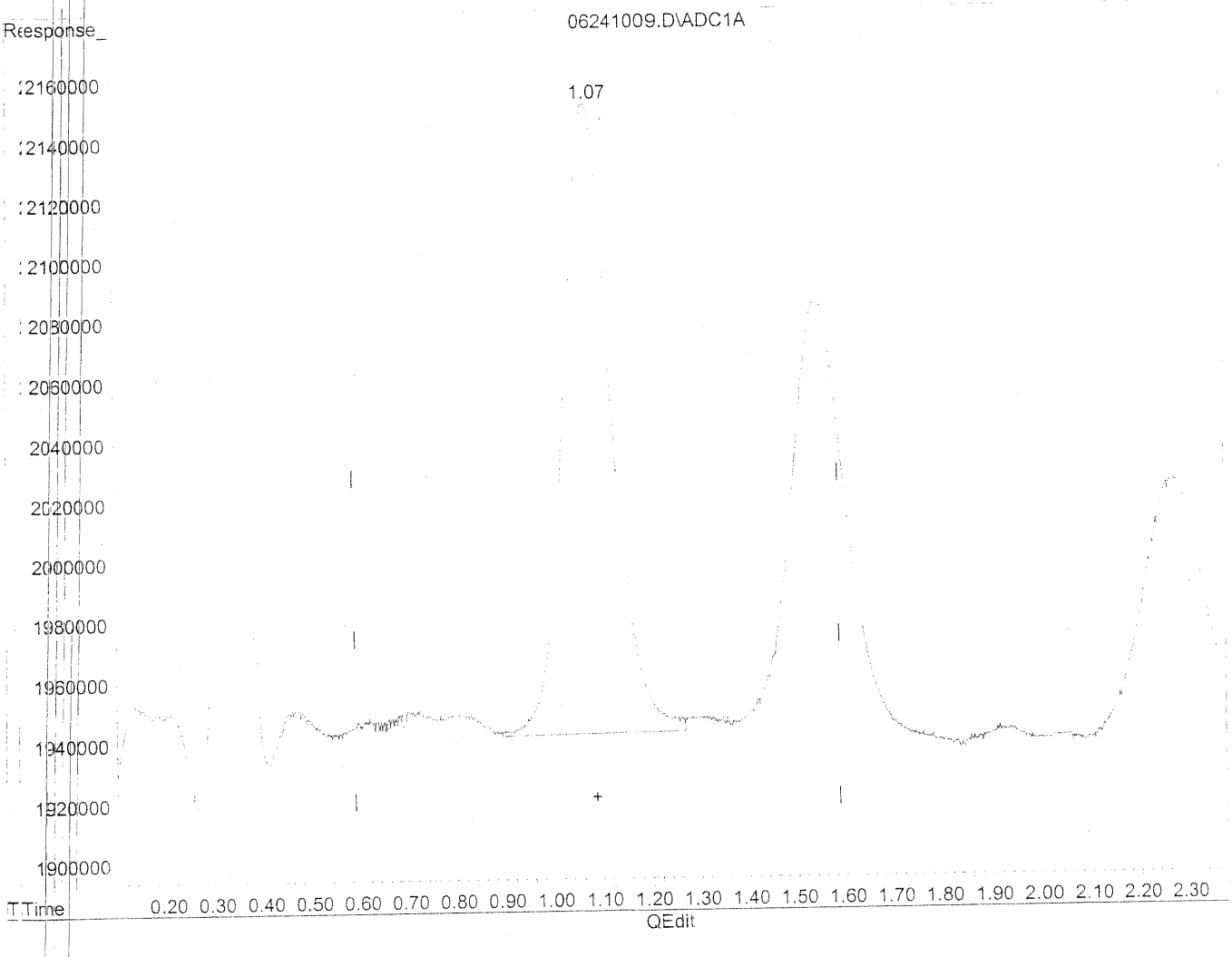
(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D
Acq On : 24 Jun 2010 13:12
Sample : 100ng/ml TO-11A S21-03091009
Misc :
IntFile : autoint1.e
Quant Time: Jun 24 13:41 19110 Quant Results File: TO110610.RES

Vial: 7
Operator: MD
Inst : LC 01
Multiplr: 1.00

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.07min 74.751ng/ml m

response 14186865

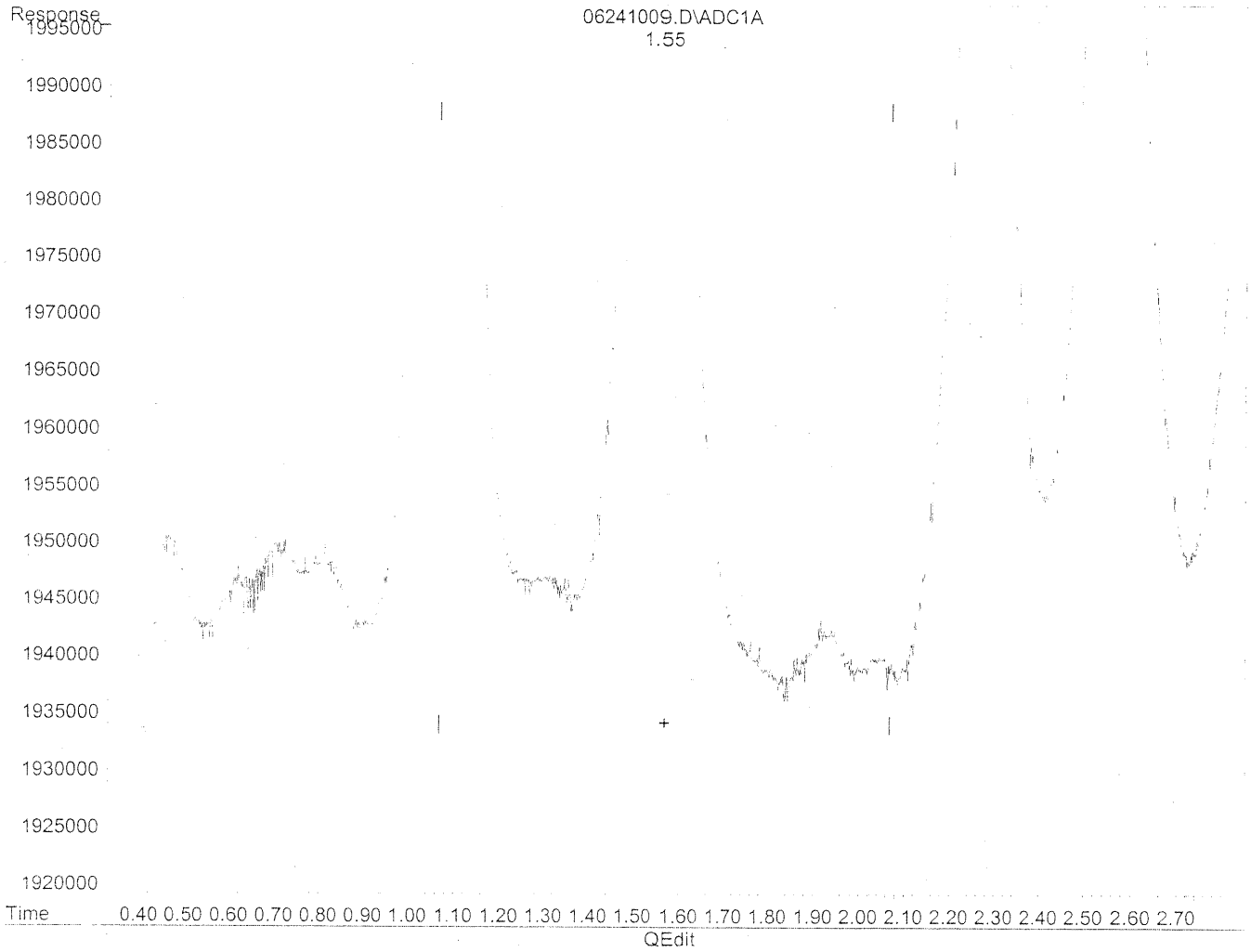
157

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:44 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.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



(2) Acetaldehyde

1.55min 82.644ng/ml

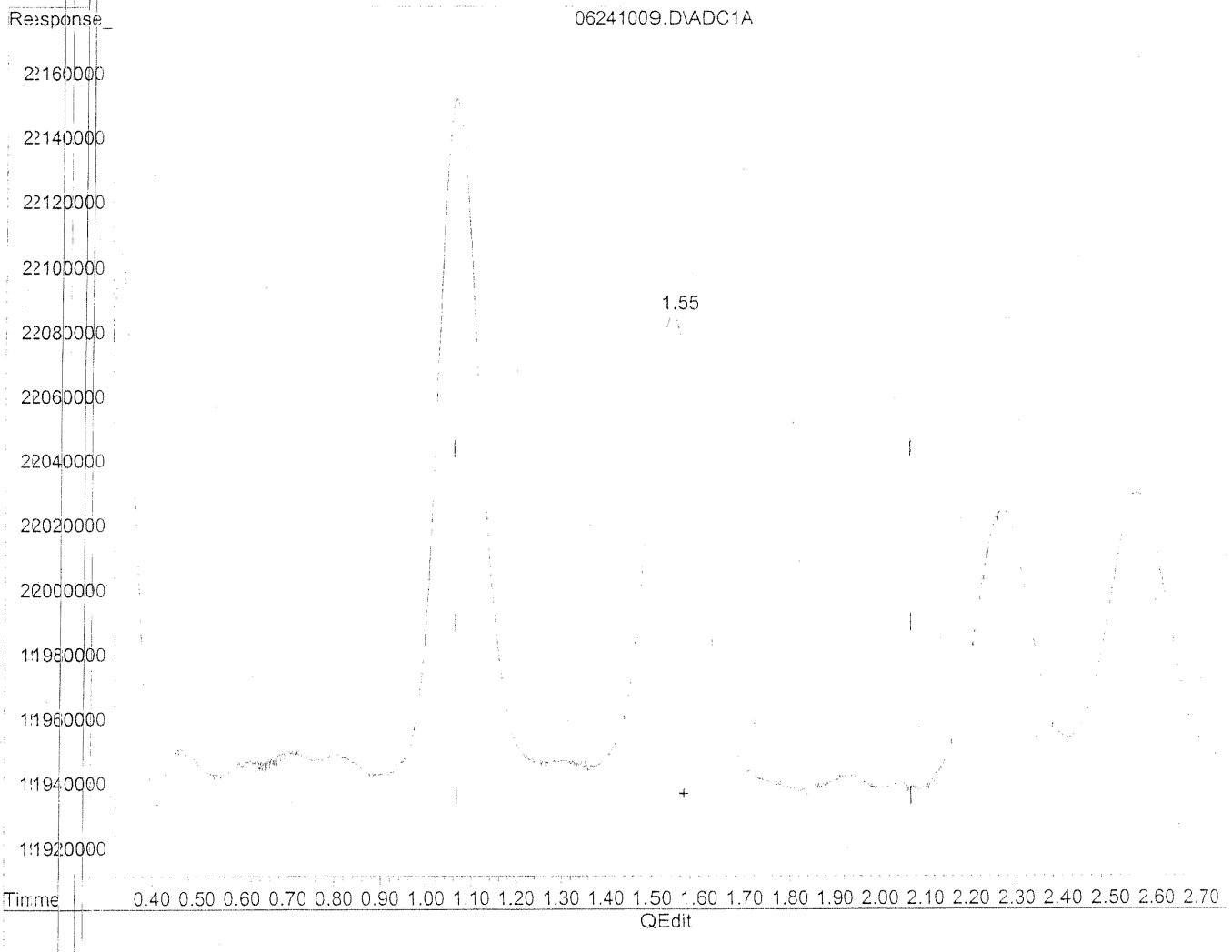
response 12079982

(+) = Expected Retention Time
06241009.D TO110610.M Thu Jun 24 16:14:17 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:47 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
Response via : Multiple Level Calibration



(2) Acetaldehyde

1.55min 76.129ng/ml m

response 11127714

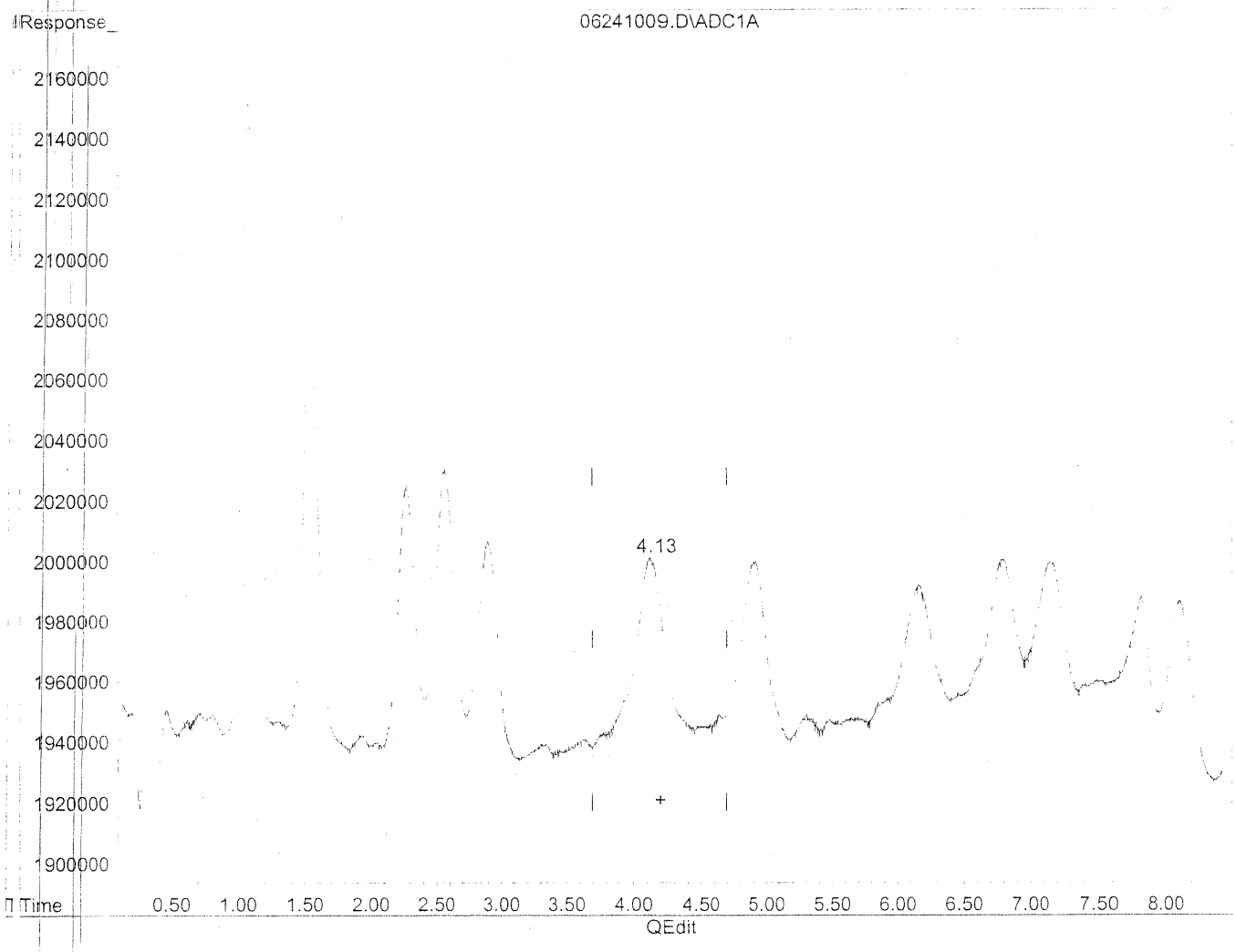
159

((+) = Expected Retention Time
06241009.D TO110610.M Fri Jun 25 10:25:37 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:41 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(4) Crotonaldehyde

4.13min 115.581ng/ml

response 11492514

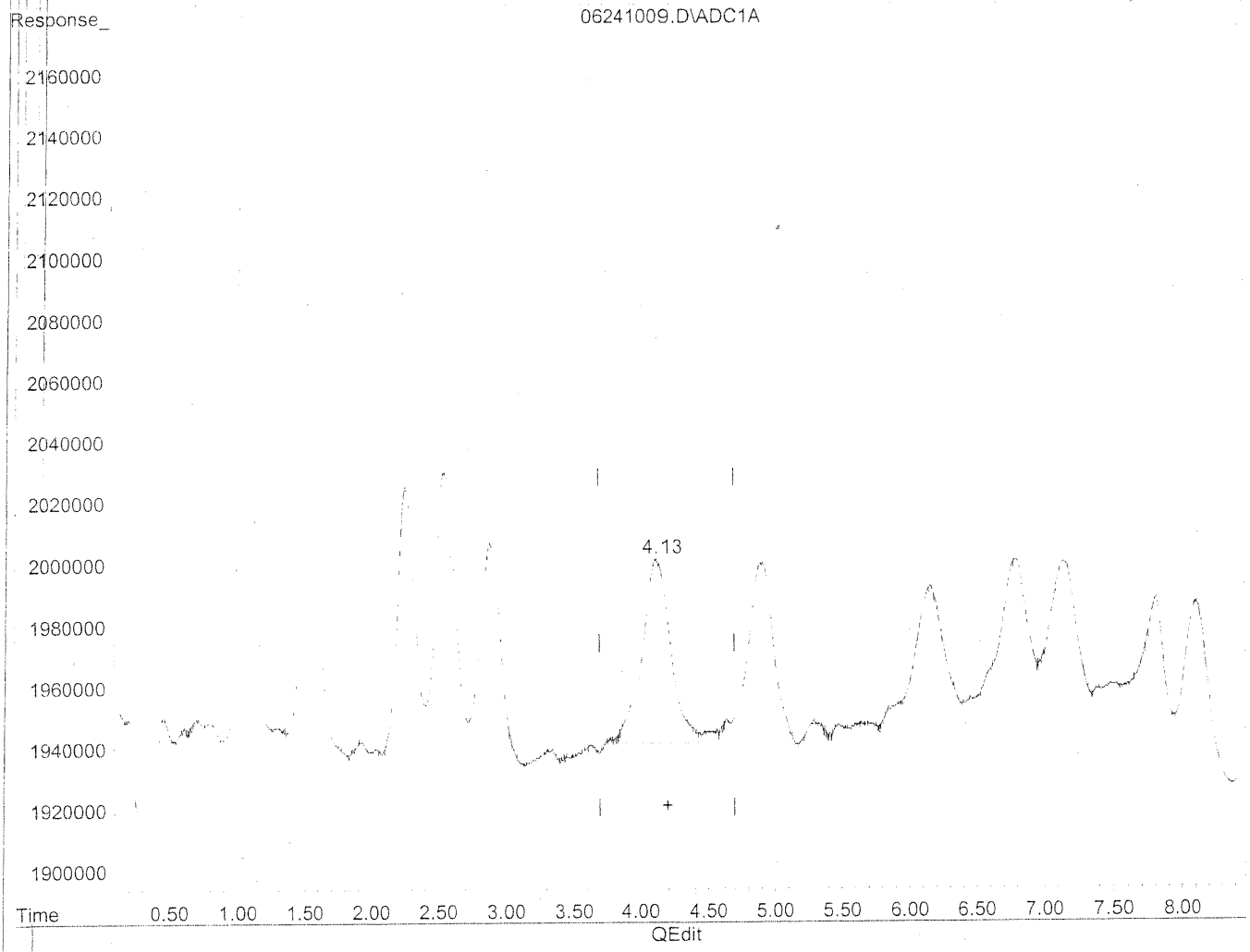
160

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:44 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.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



(4) Crotonaldehyde

4.13min 85.389ng/ml m

response 8490383

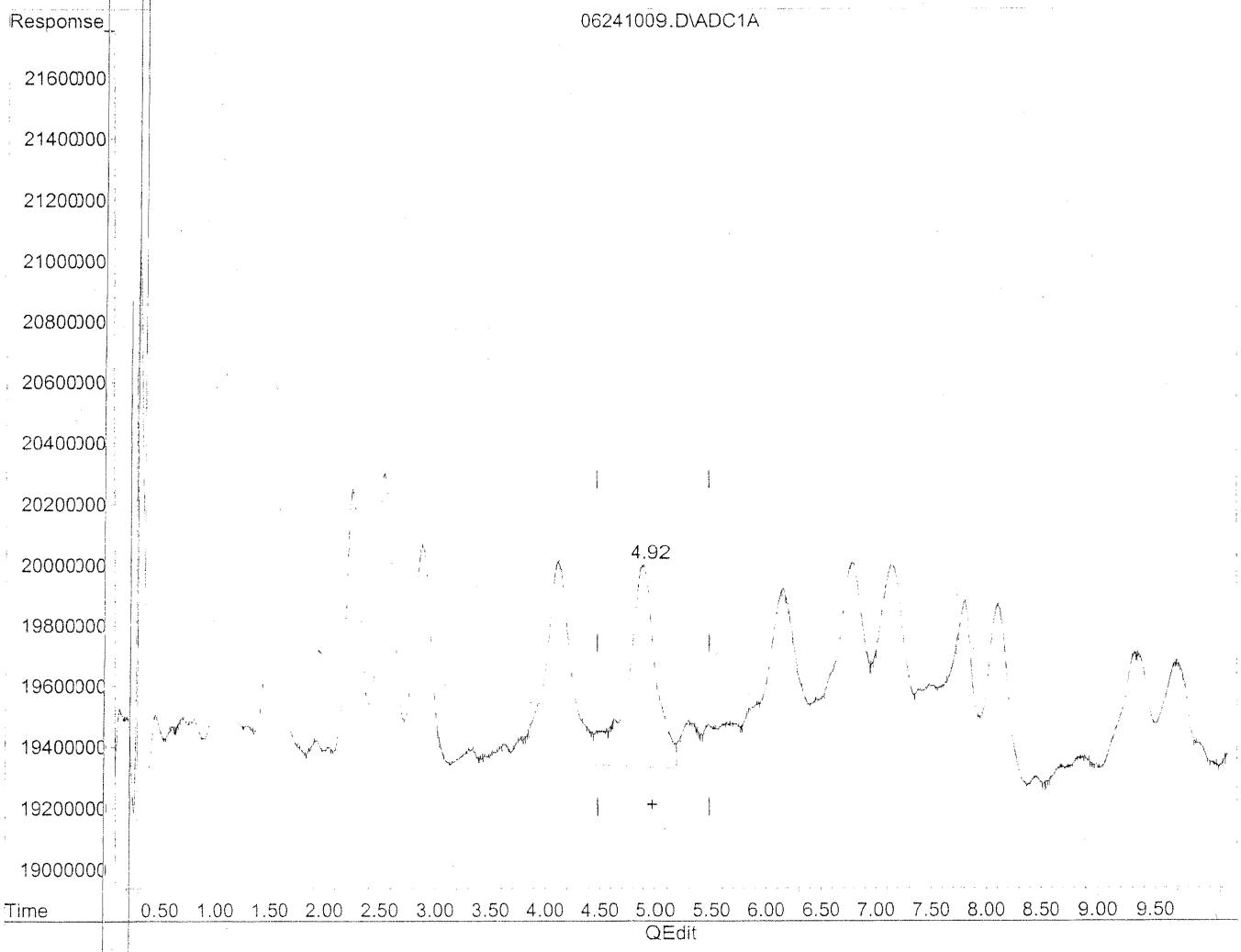
161

(+) = Expected Retention Time
06241009.D TO110610.M Thu Jun 24 16:22:18 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:41 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(5) Butyraldehyde

4.91min 127.486ng/ml

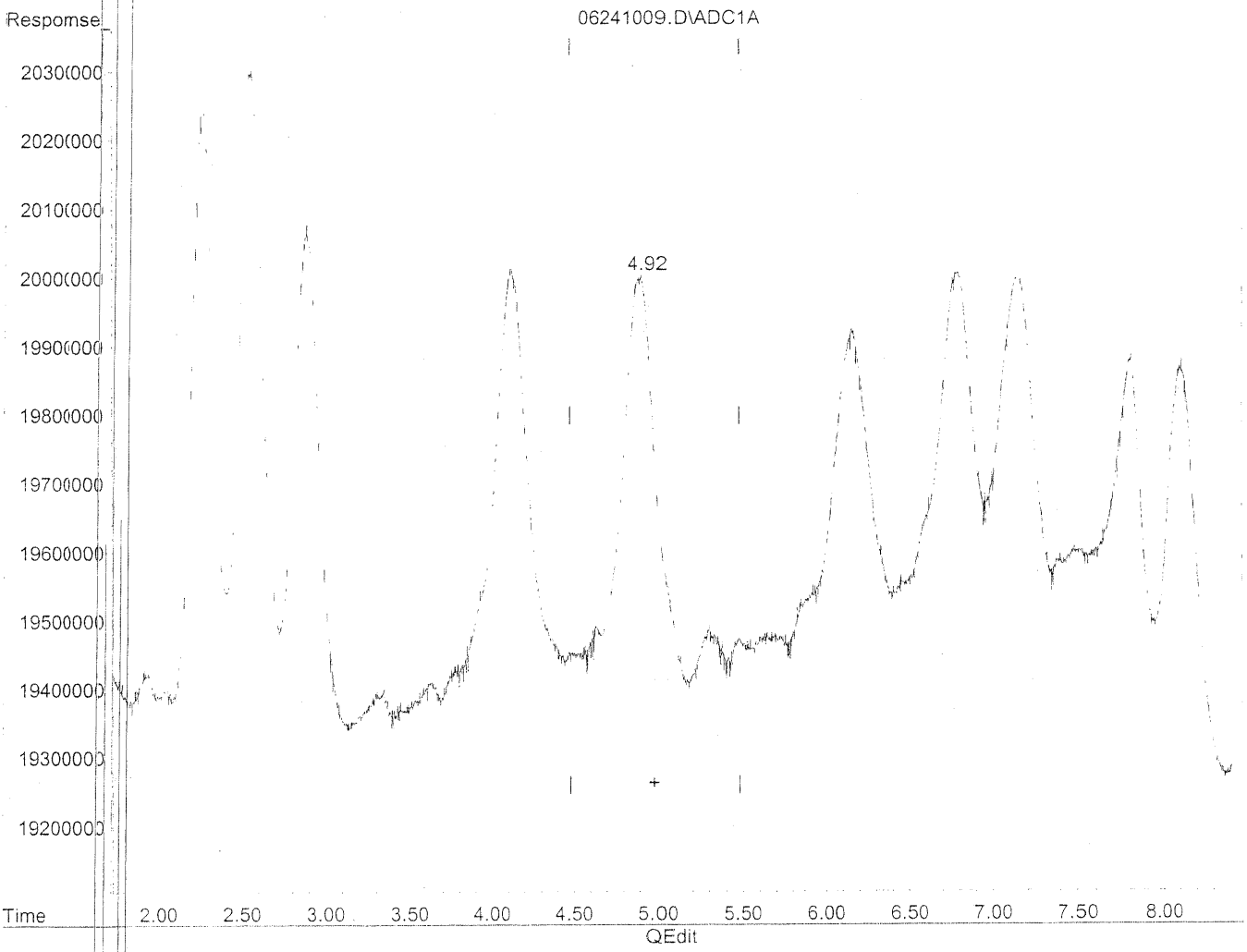
response 11898098

(+)) = Expected Retention Time
062441009.D TO110610.M Thu Jun 24 13:42:05 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:44 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.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



(5) Butyraldehyde

4.92min 86.753ng/ml m

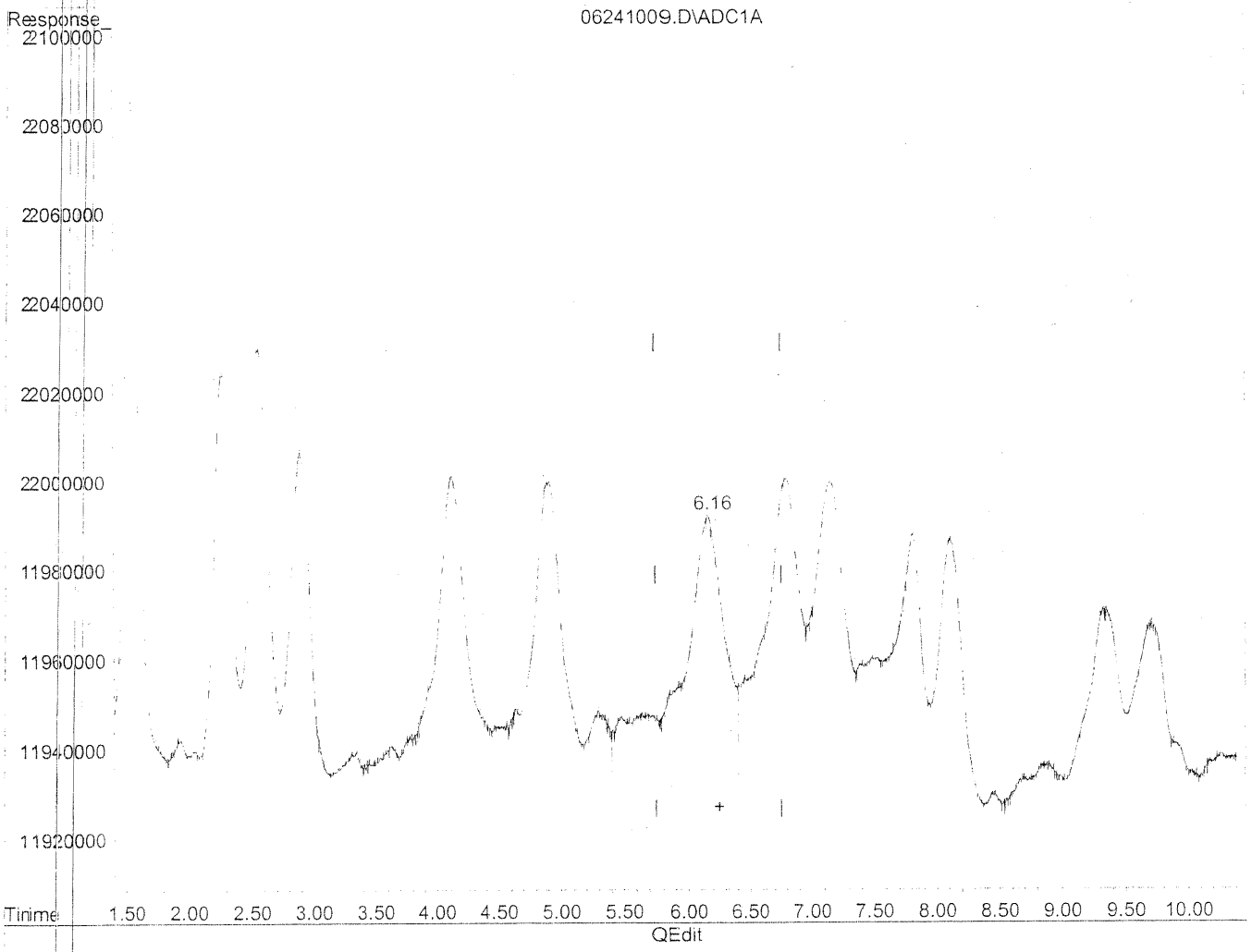
response 8096555

(+)) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:41 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.16min 232.713ng/ml

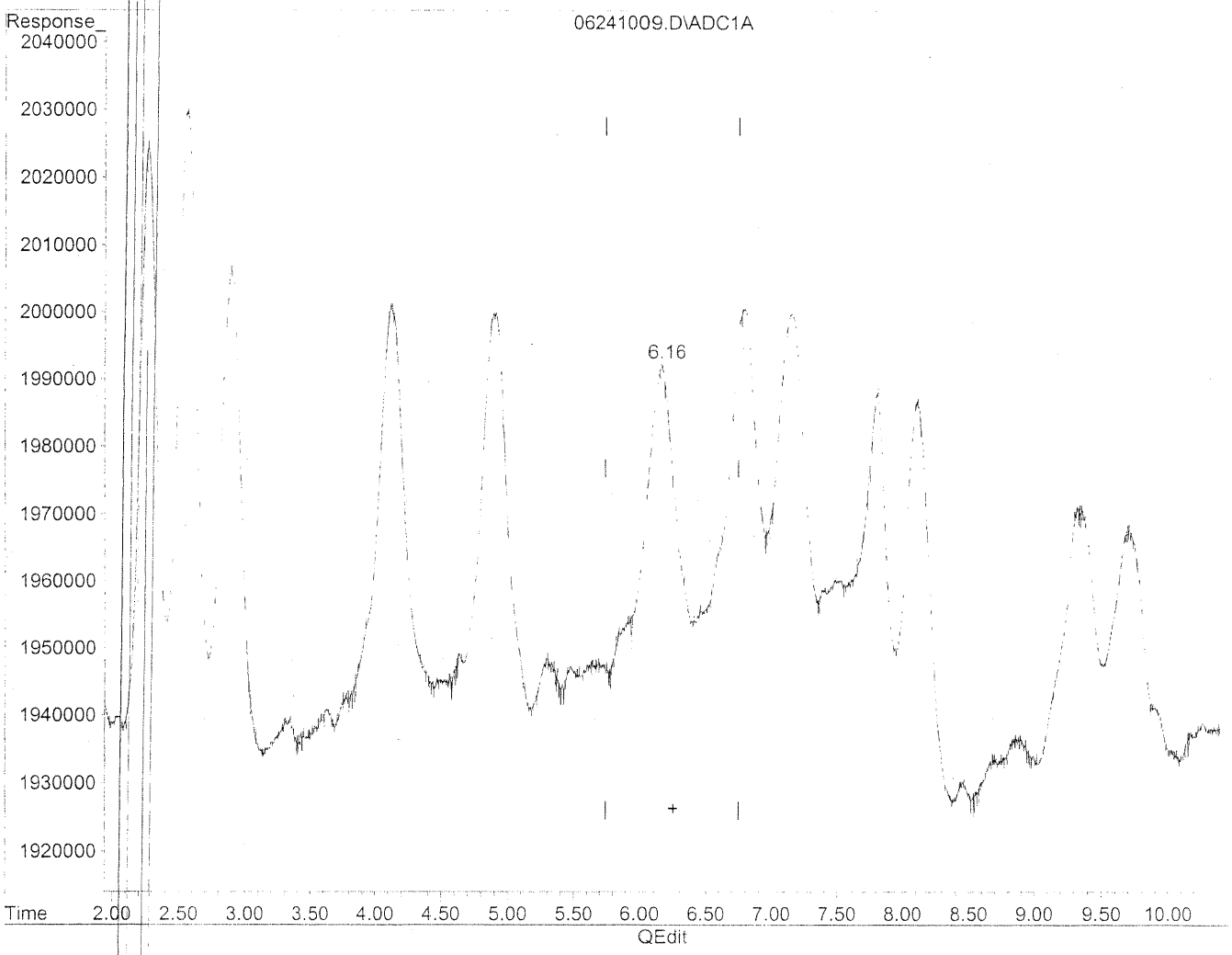
response 16196772

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:47 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
Response via : Multiple Level Calibration



(6) Benzaldehyde

6.16min 84.678ng/ml m

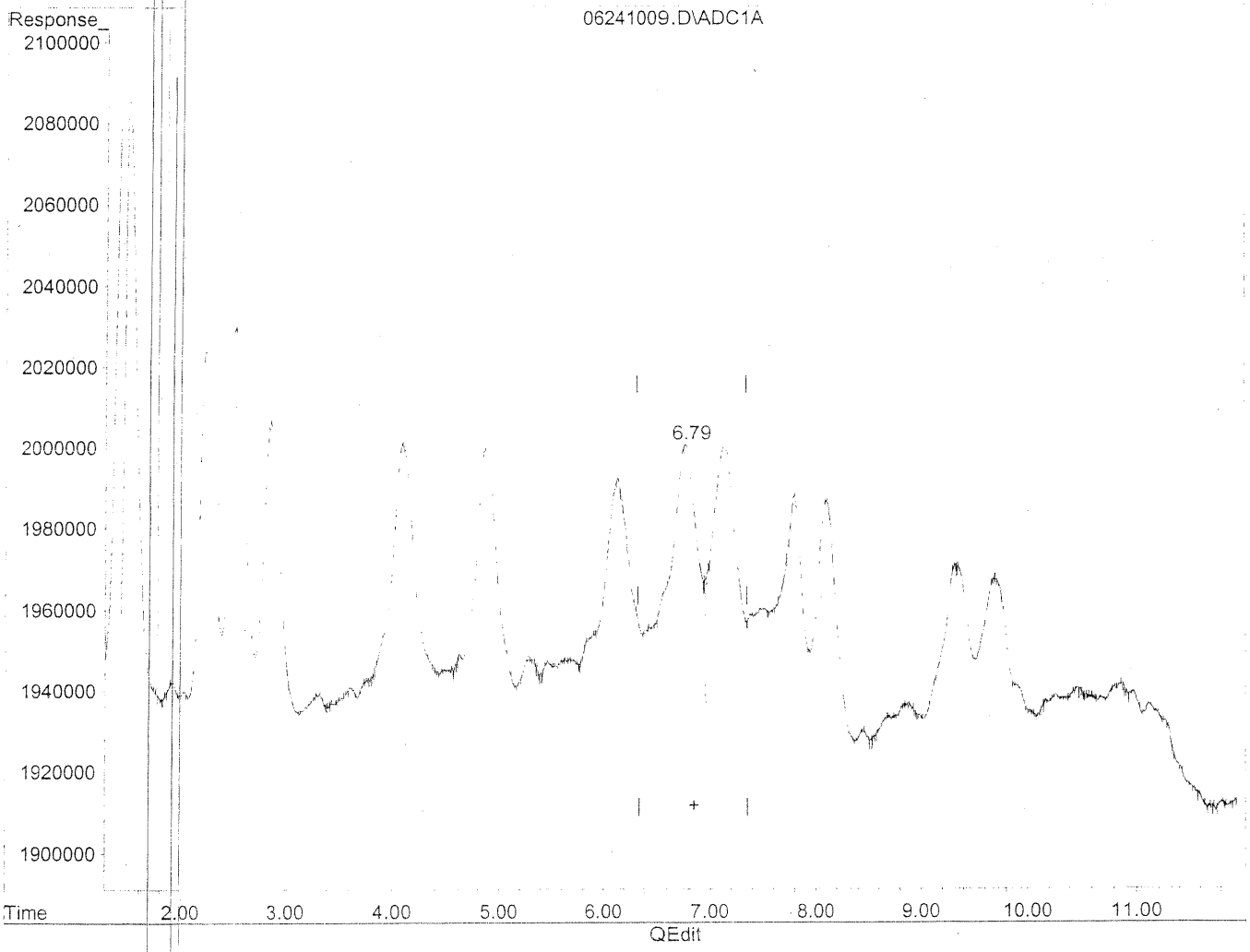
response 5893577

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:41 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(7) Isovaleraldehyde

6 6.79min 177.899ng/ml

r response 14477157

166

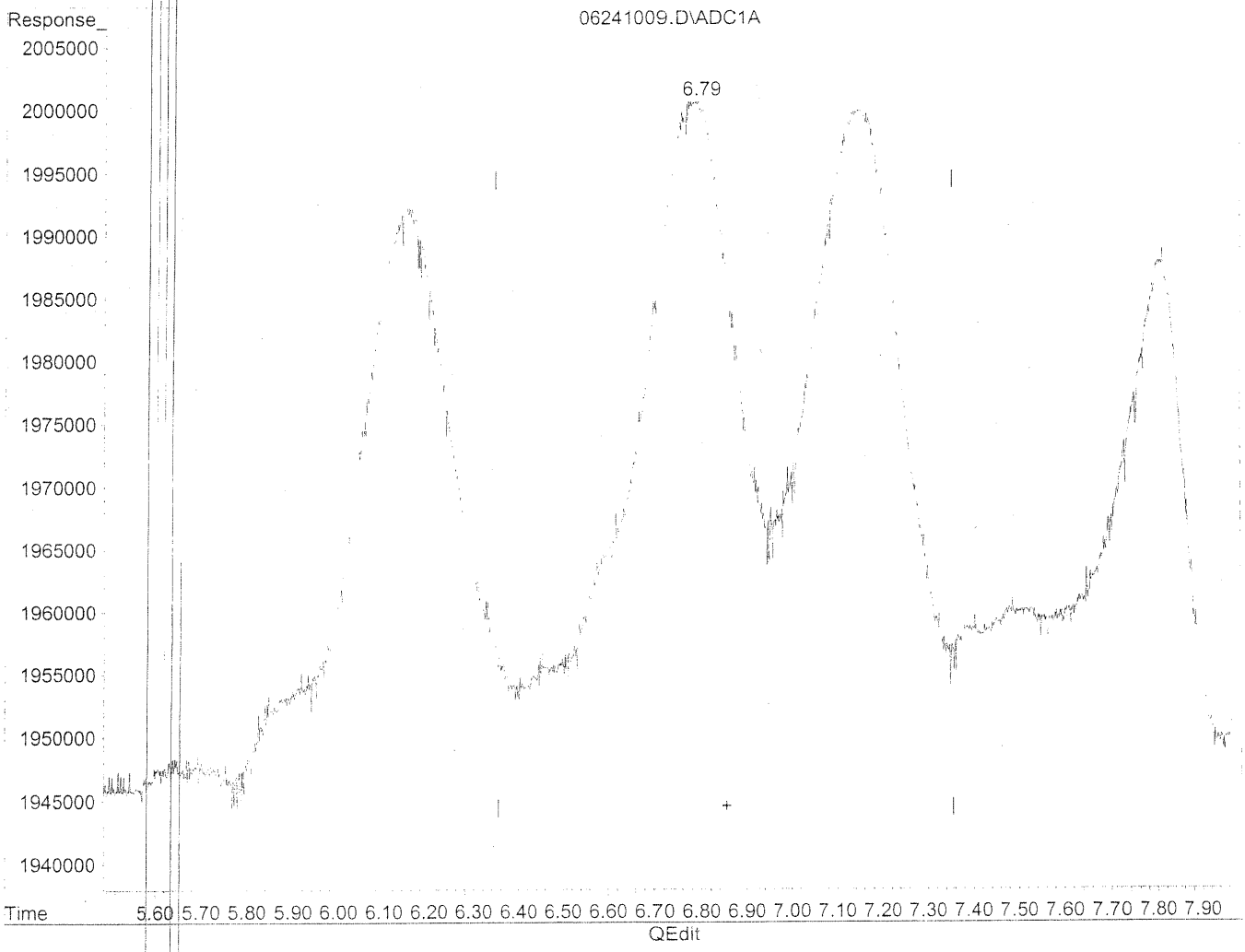
(+) = Expected Retention Time

062410009.D TO110610.M Thu Jun 24 13:42:55 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:24 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.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

6 6.79min 82.699ng/ml m

response 6729977

167

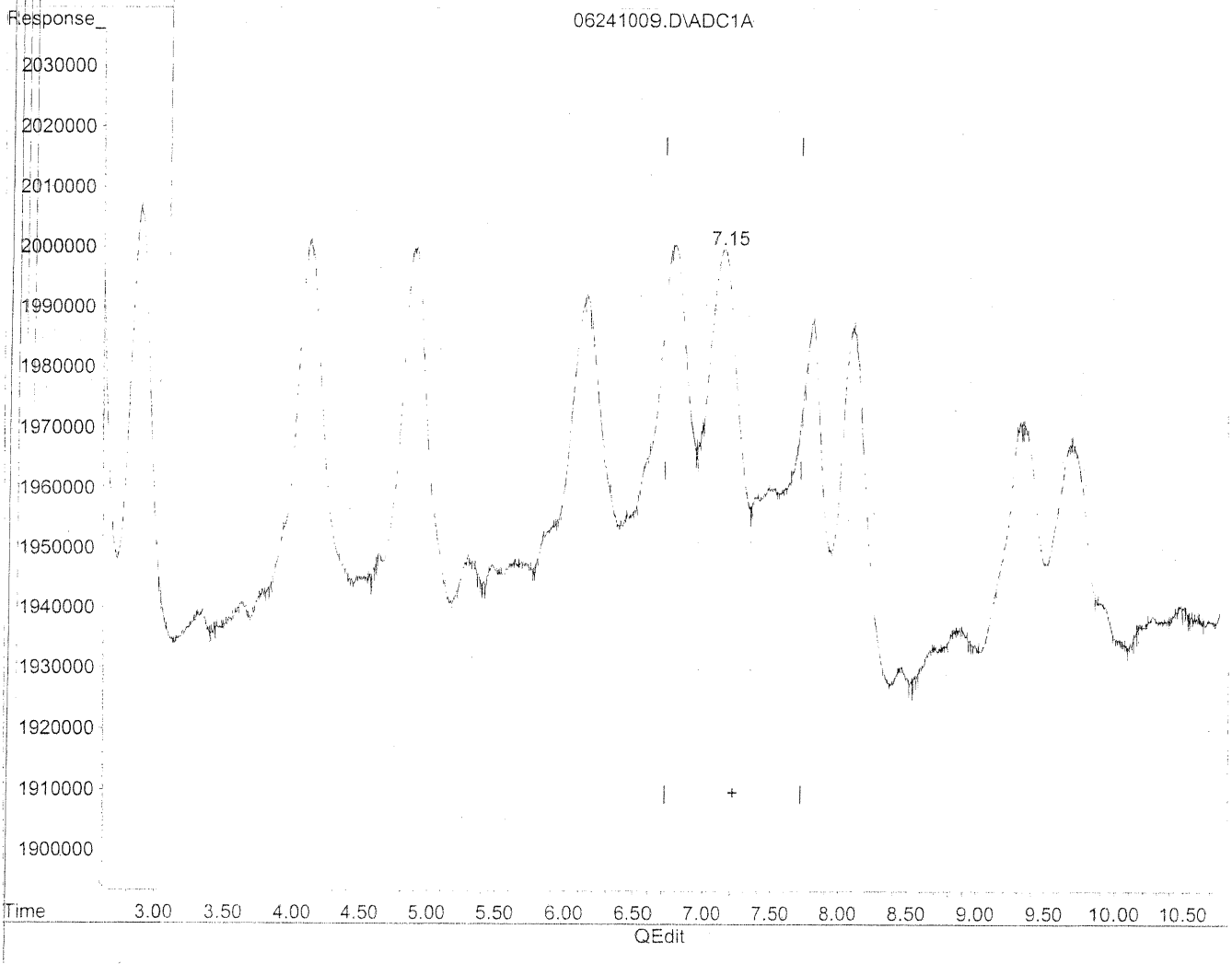
(+) = Expected Retention Time

062410009.D TO110610.M Thu Jun 24 16:26:28 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:41 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.15min 157.780ng/ml

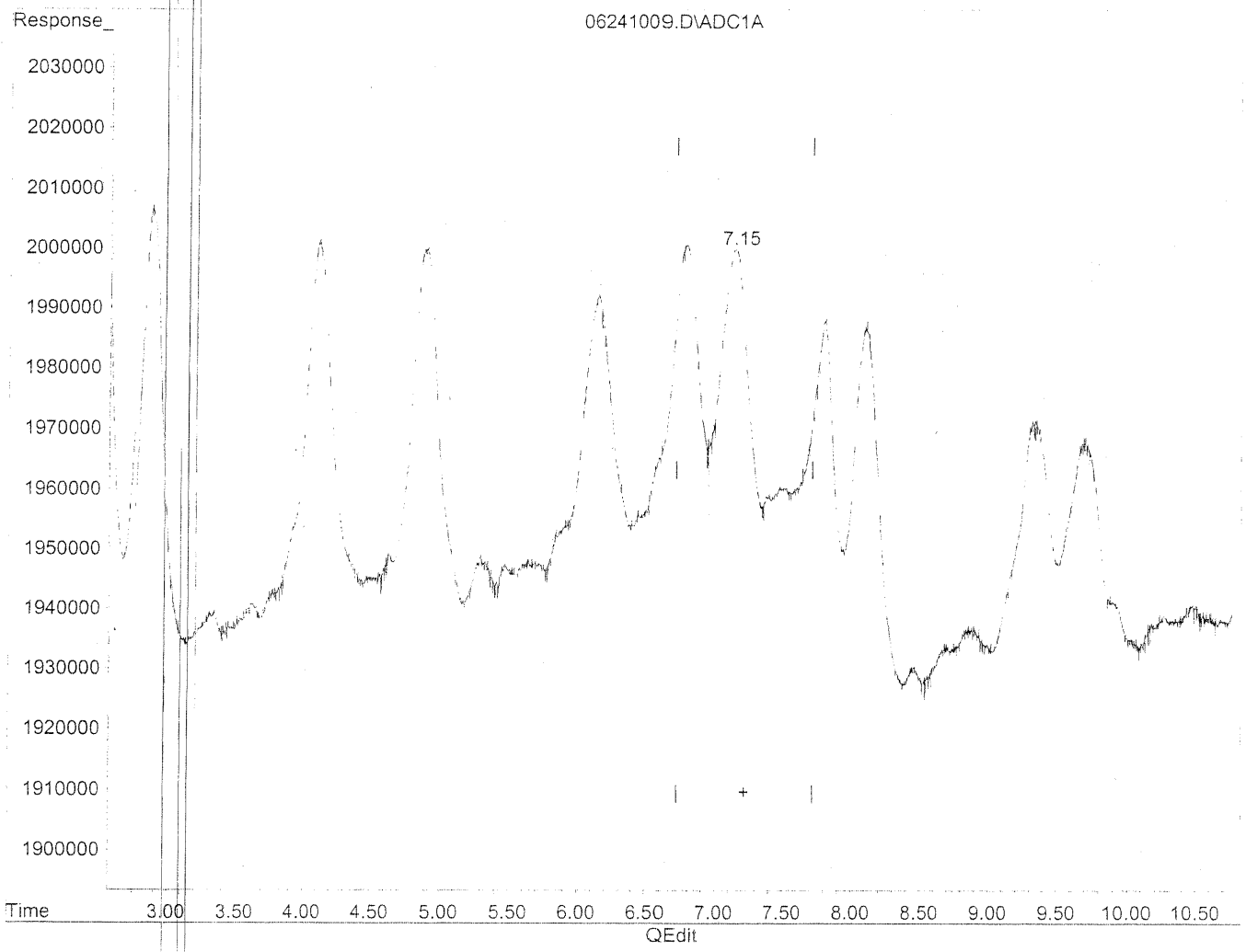
response 11853035

(+) = Expected Retention Time
06241009.D TO110610.M Thu Jun 24 13:43:20 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:41 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(8) Valeraldehyde

7.15min 88.470ng/ml m

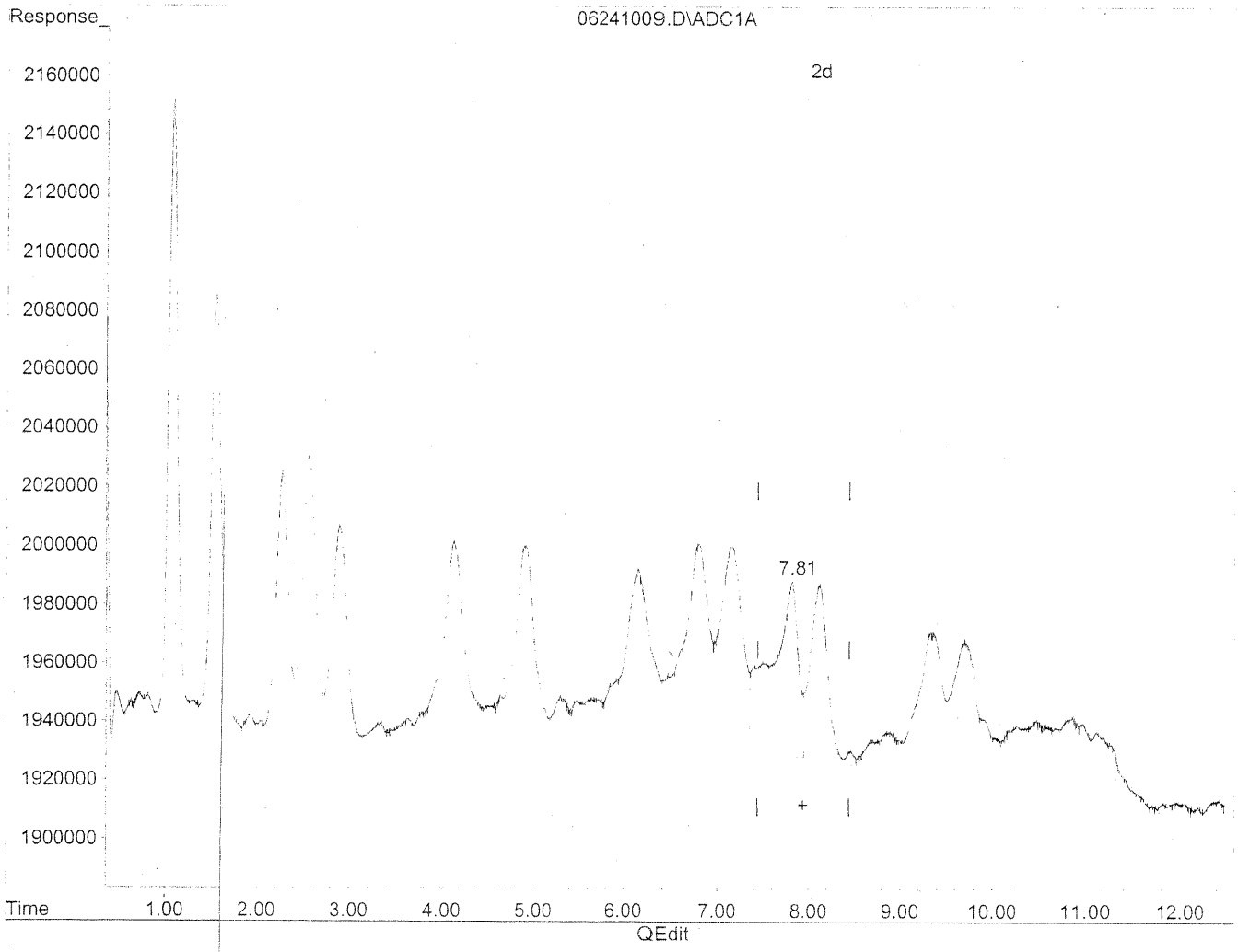
response 6646232

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:41 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

7.81min 202.040ng/ml

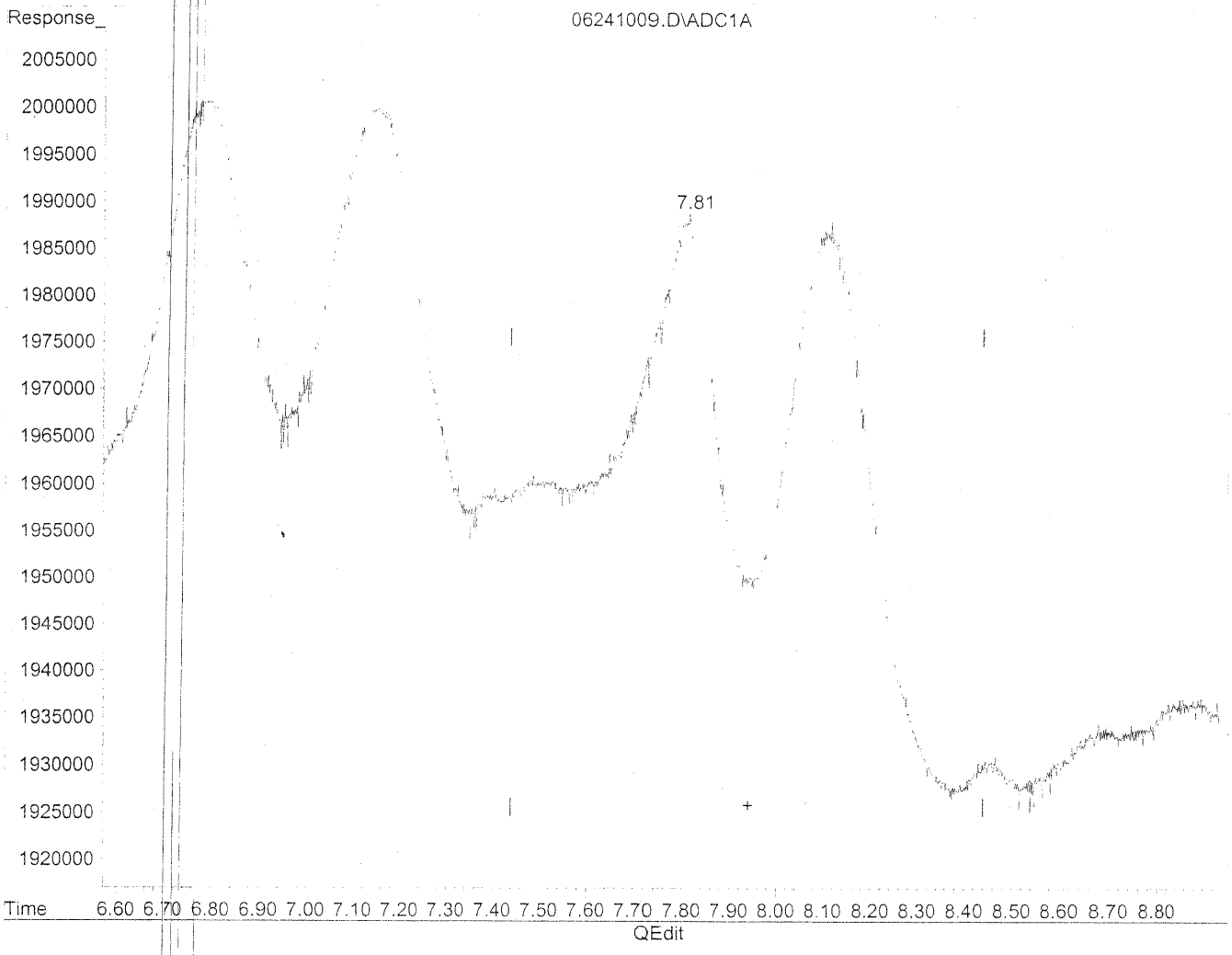
response 12862545

(+) = Expected Retention Time
06241009.D TO110610.M Thu Jun 24 13:43:40 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:47 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 16:50:51 2010
Response via : Multiple Level Calibration



(9) o-Tolualdehyde

7.81min 71.724ng/ml m

response 4566176

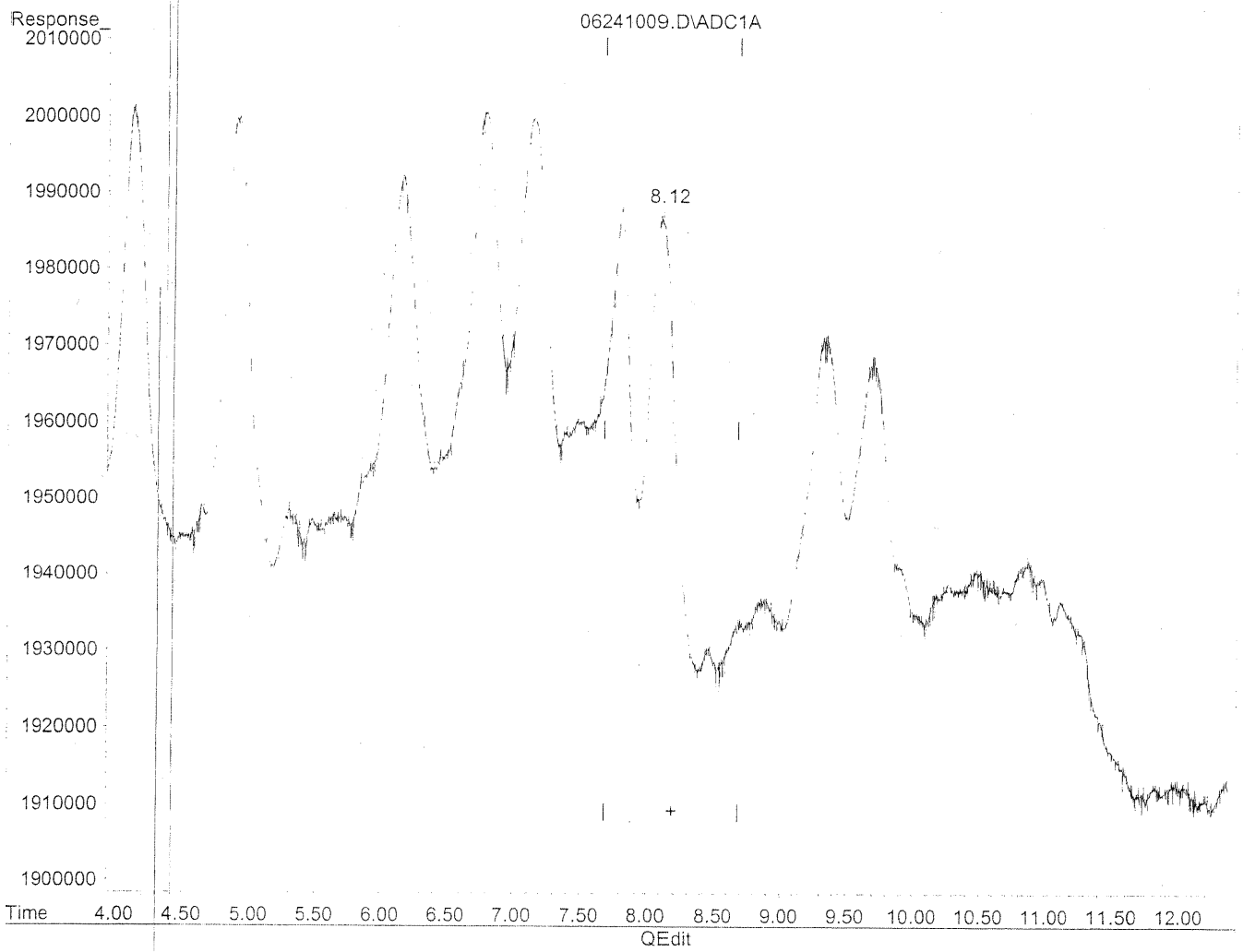
171

(+) = Expected Retention Time
06241009.D TO110610.M Thu Jun 24 16:51:02 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:44 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.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



(10) m,p-Tolualdehyde

8.11min 133.544ng/ml

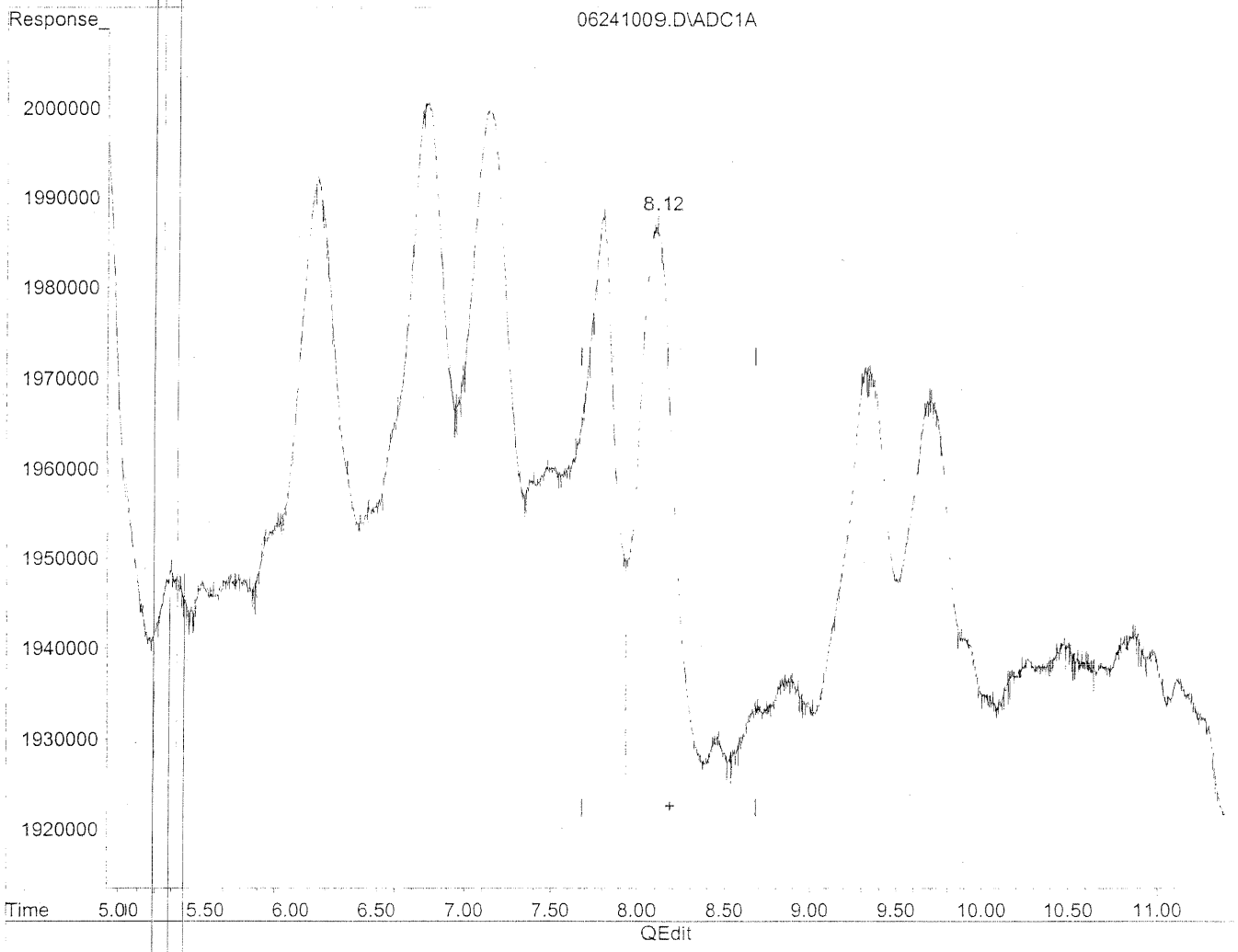
response 7572174

(+) = Expected Retention Time
06241009.D TO110610.M Thu Jun 24 16:24:29 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:47 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

8.12min 143.539ng/ml m

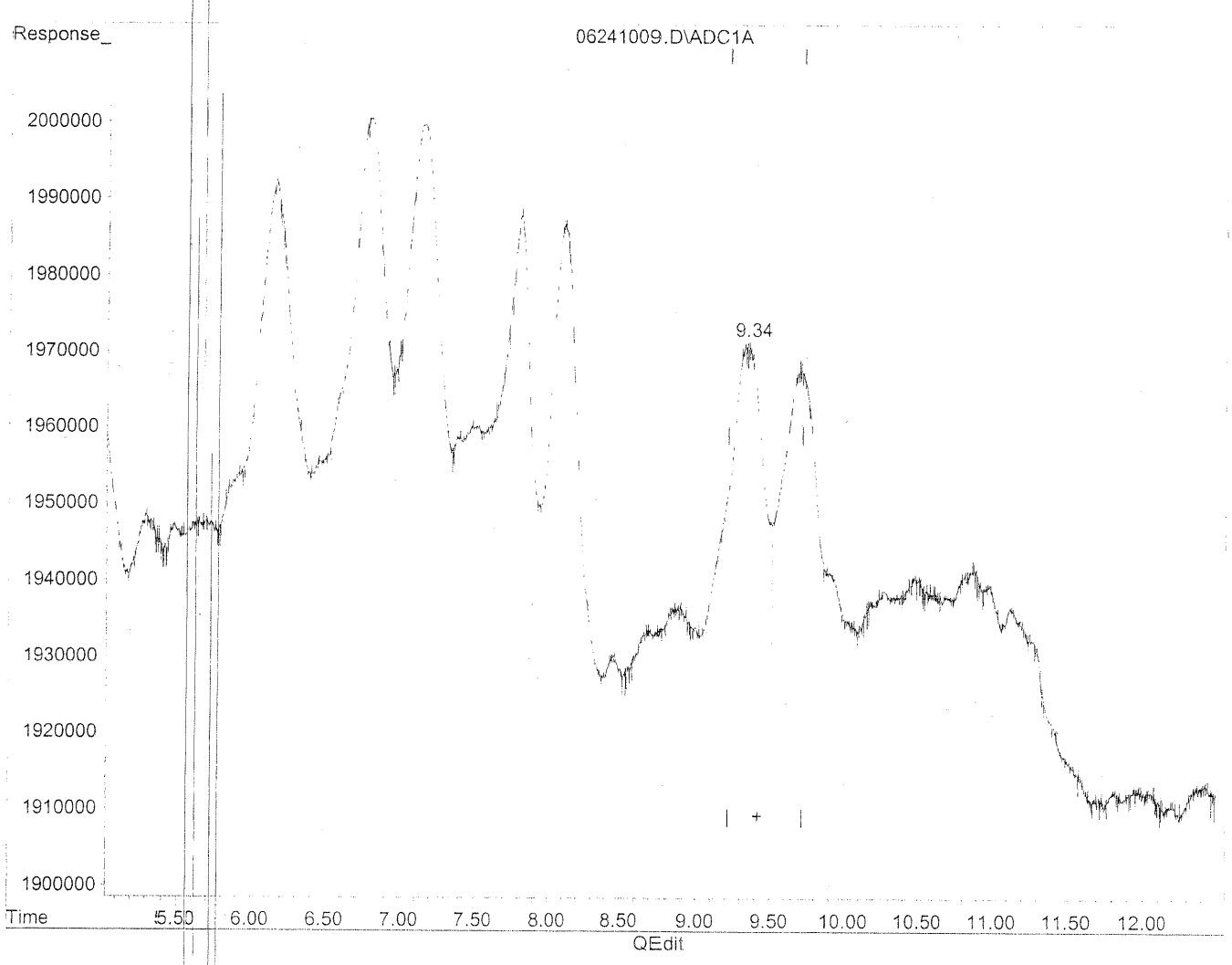
response 8138944

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:41 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(11) Hexaldehyde

9.33min 94.666ng/ml

response 6757167

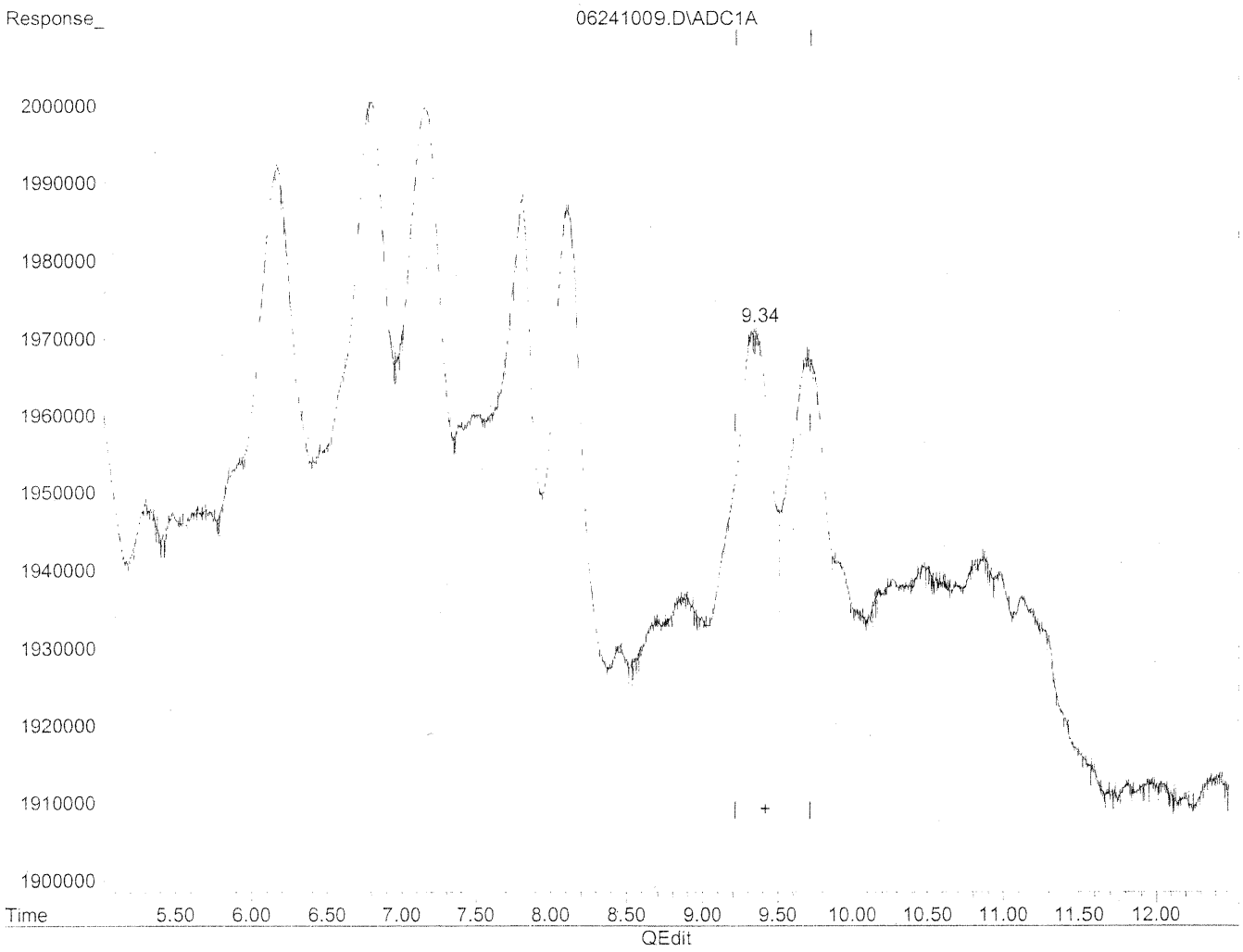
174

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:44 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.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



(11) Hexaldehyde

9.34min 82.359ng/ml m

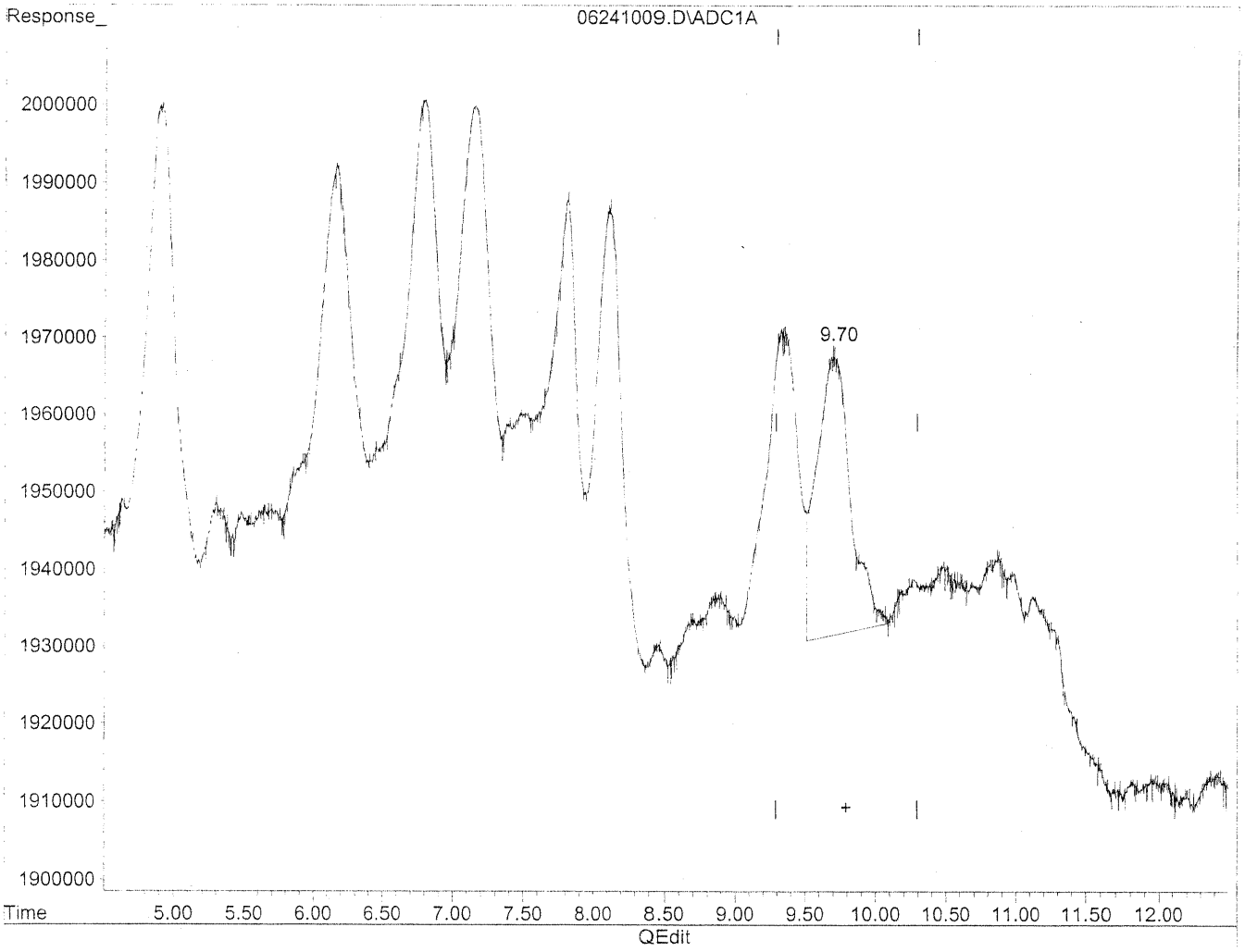
response 5878688

(+) = Expected Retention Time
06241009.D TO110610.M Thu Jun 24 16:25:12 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:41 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

9.70min 123.594ng/ml

response 6071575

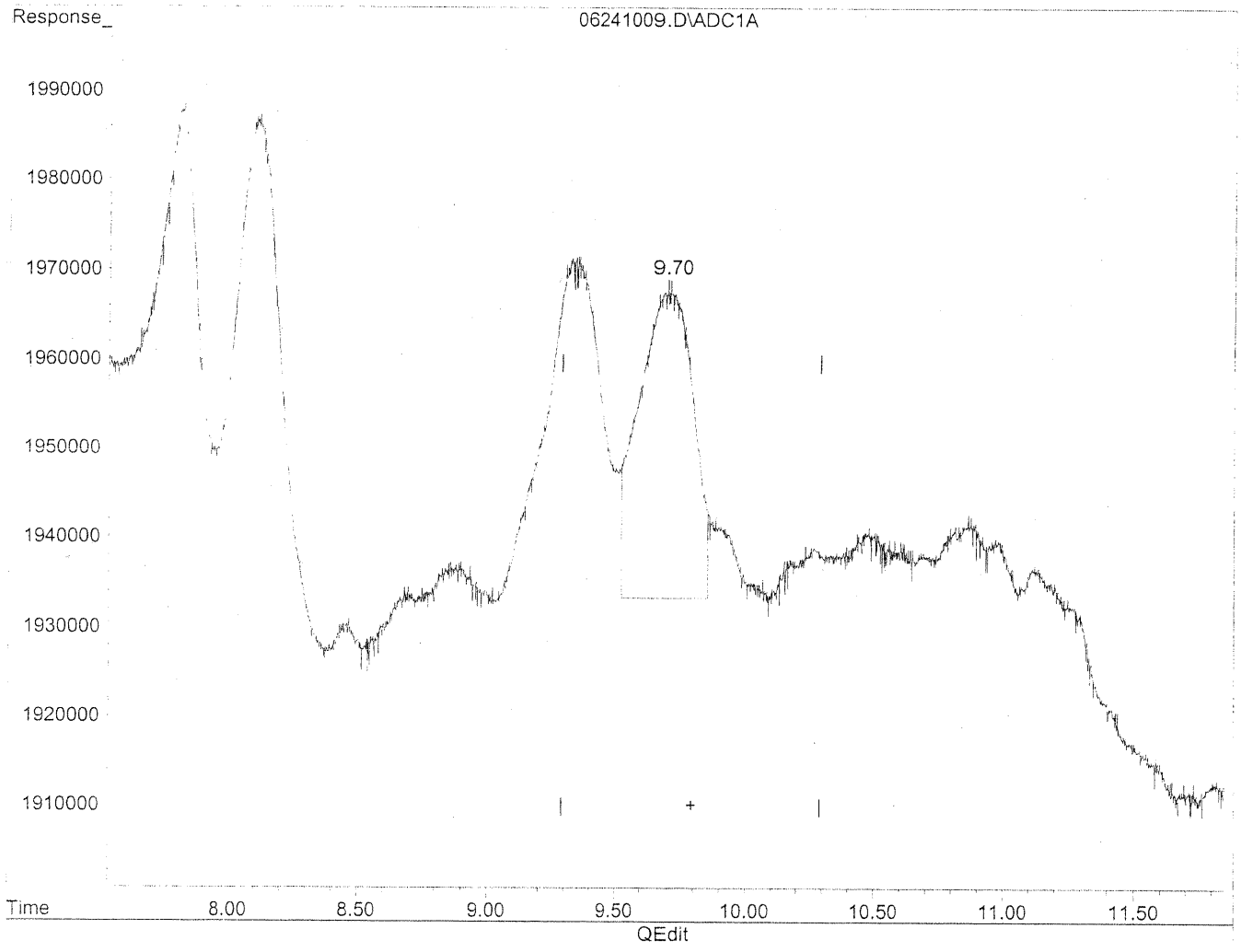
176

(+) = Expected Retention Time
06241009.D TO110610.M Thu Jun 24 13:44:29 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241009.D Vial: 7
Acq On : 24 Jun 2010 13:12 Operator: MD
Sample : 100ng/ml TO-11A S21-03091009 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:47 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 16:50:51 2010
Response via : Multiple Level Calibration



(12) 2,5-Dimethylbenzaldehyde

9.70min 98.946ng/ml m

response 4860735

5/3/10

4/26/10

177

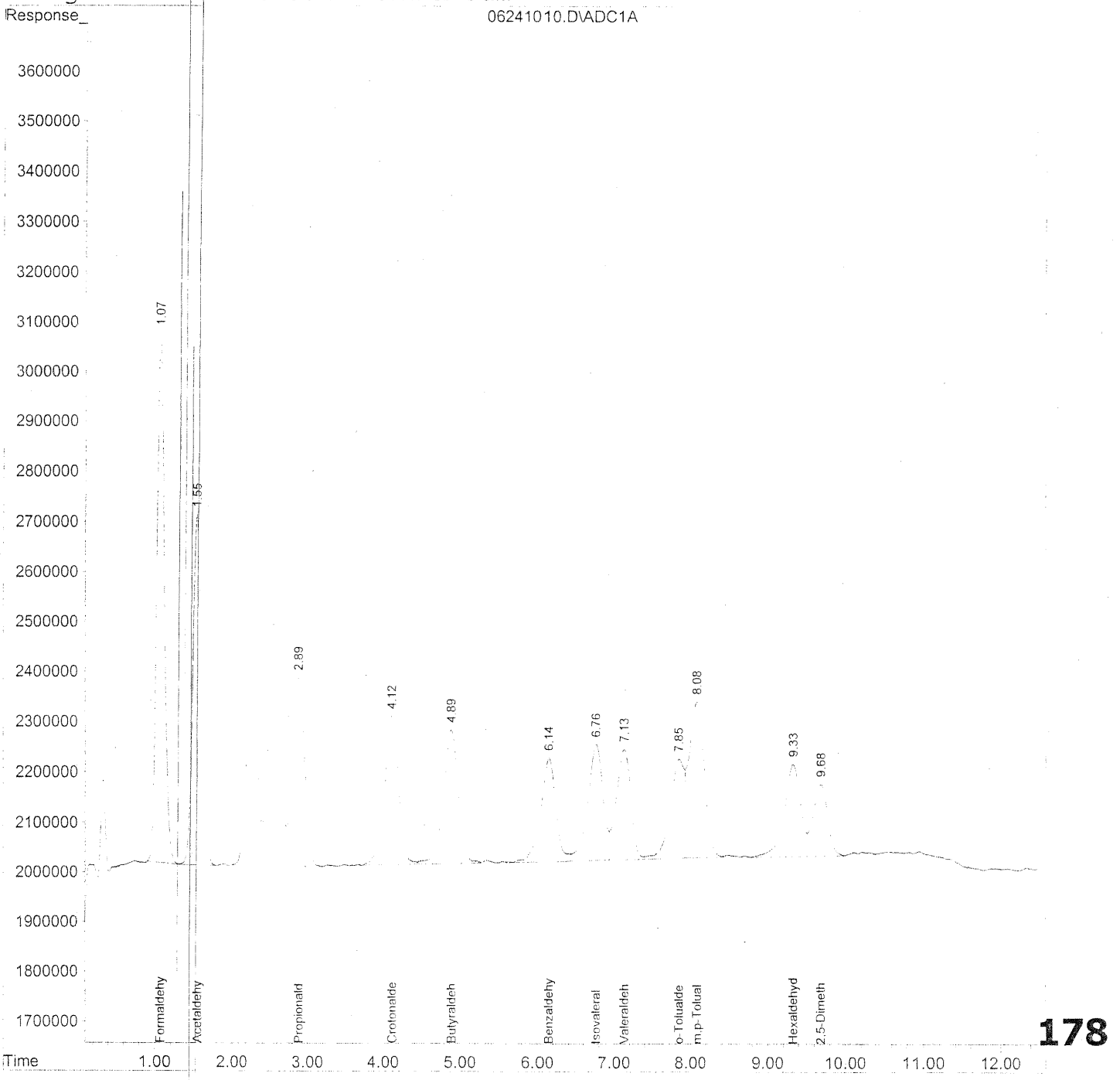
(+) = Expected Retention Time
06241009.D TO110610.M Thu Jun 24 16:51:12 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241010.D Vial: 8
Acq On : 24 Jun 2010 13:25 Operator: MD
Sample : 500ng/ml TO-11A S21-03091008 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:45 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
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\2010_06\24\06241010.D Vial: 8
Acq On : 24 Jun 2010 13:25 Operator: MD
Sample : 500ng/ml TO-11A S21-03091008 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:45 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:26:29 2010
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	1.07	71248470	374.763	ng/ml
2)	Acetaldehyde	1.55	54504441	370.871	ng/ml
3)	Propionaldehyde	2.88	41248067	375.715	ng/ml
4)	Crotonaldehyde	4.12	38090464	379.879	ng/ml
5)	Butyraldehyde	4.89	35410354	376.212	ng/ml
6)	Benzaldehyde	6.14	28344164	408.486	ng/ml
7)	Isovaleraldehyde	6.76	32418035	399.746	ng/ml
8)	Valeraldehyde	7.13	28548123	378.106	ng/ml
9)	o-Tolualdehyde	7.85	24943255	395.514	ng/ml
10)	m,p-Tolualdehyde	8.08	43217145	767.147	ng/ml
11)	Hexaldehyde	9.33	27064970	377.481	ng/ml
12)	2,5-Dimethylbenzaldehyde	9.68	18911052	380.246	ng/ml

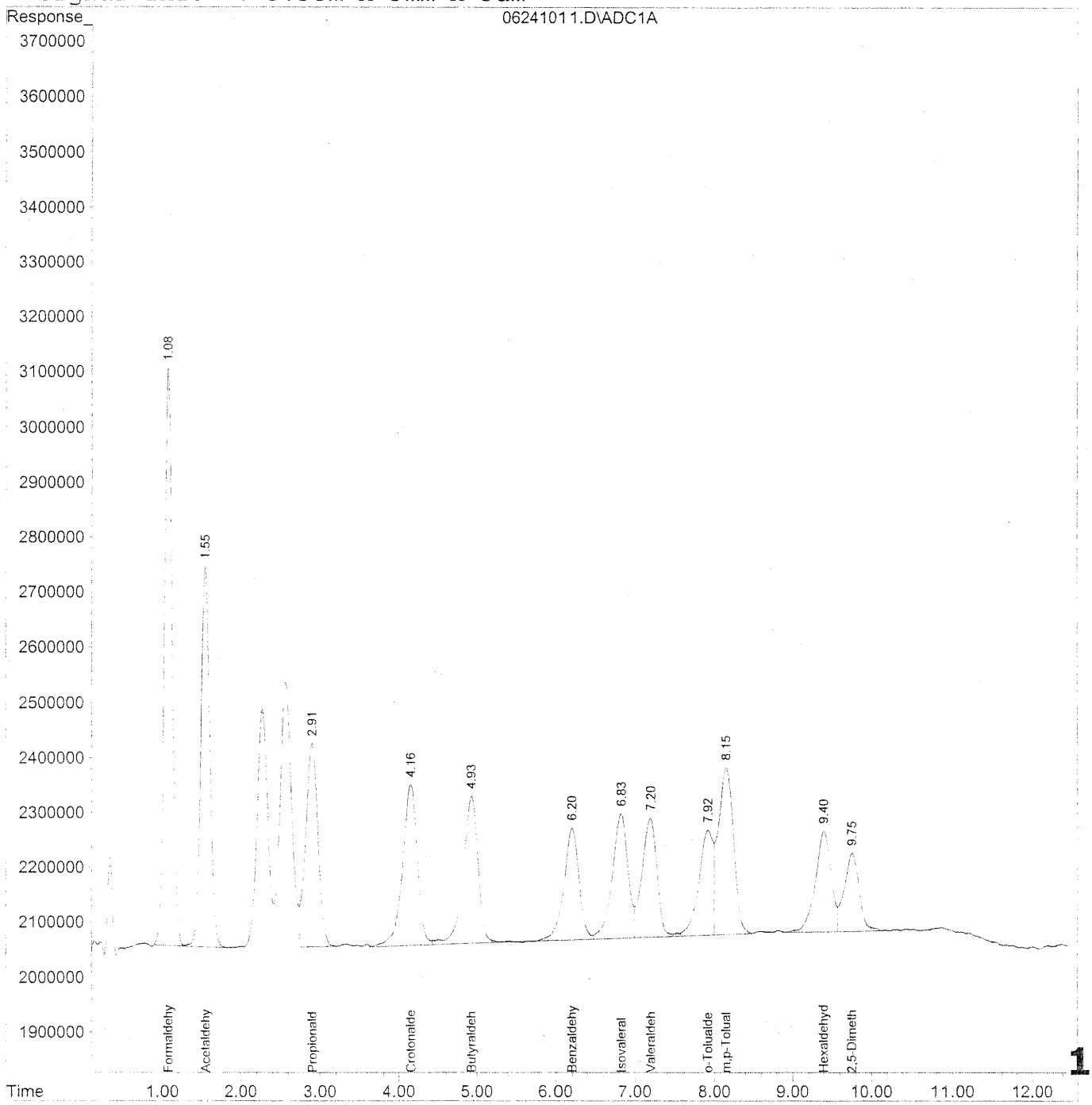
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241011.D Vial: 9
Acq On : 24 Jun 2010 13:39 Operator: MD
Sample : 500ng/ml TO-11A S21-03091008 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:55 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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\2010_06\24\06241011.D Vial: 9
Acq On : 24 Jun 2010 13:39 Operator: MD
Sample : 500ng/ml TO-11A S21-03091008 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 13:55 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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	1.08	70508340	513.050	ng/ml
2)	Acetaldehyde	1.55	54114192	497.836	ng/ml
3)	Propionaldehyde	2.91	40426693	507.179	ng/ml
4)	Crotonaldehyde	4.15	38548382	504.619	ng/ml
5)	Butyraldehyde	4.93	35949330	491.446	ng/ml
6)	Benzaldehyde	6.20	26799052	476.450	ng/ml
7)	Isovaleraldehyde	6.83	31444882	492.141	ng/ml
8)	Valeraldehyde	7.19	28285881	489.439	ng/ml
9)	o-Tolualdehyde	7.92	23916376	493.631	ng/ml
10)	m,p-Tolualdehyde	8.15	42173453	997.424	ng/ml
11)	Hexaldehyde	9.40	25614342	478.153	ng/ml
12)	2,5-Dimethylbenzaldehyde	9.76	18988988	478.456	ng/ml

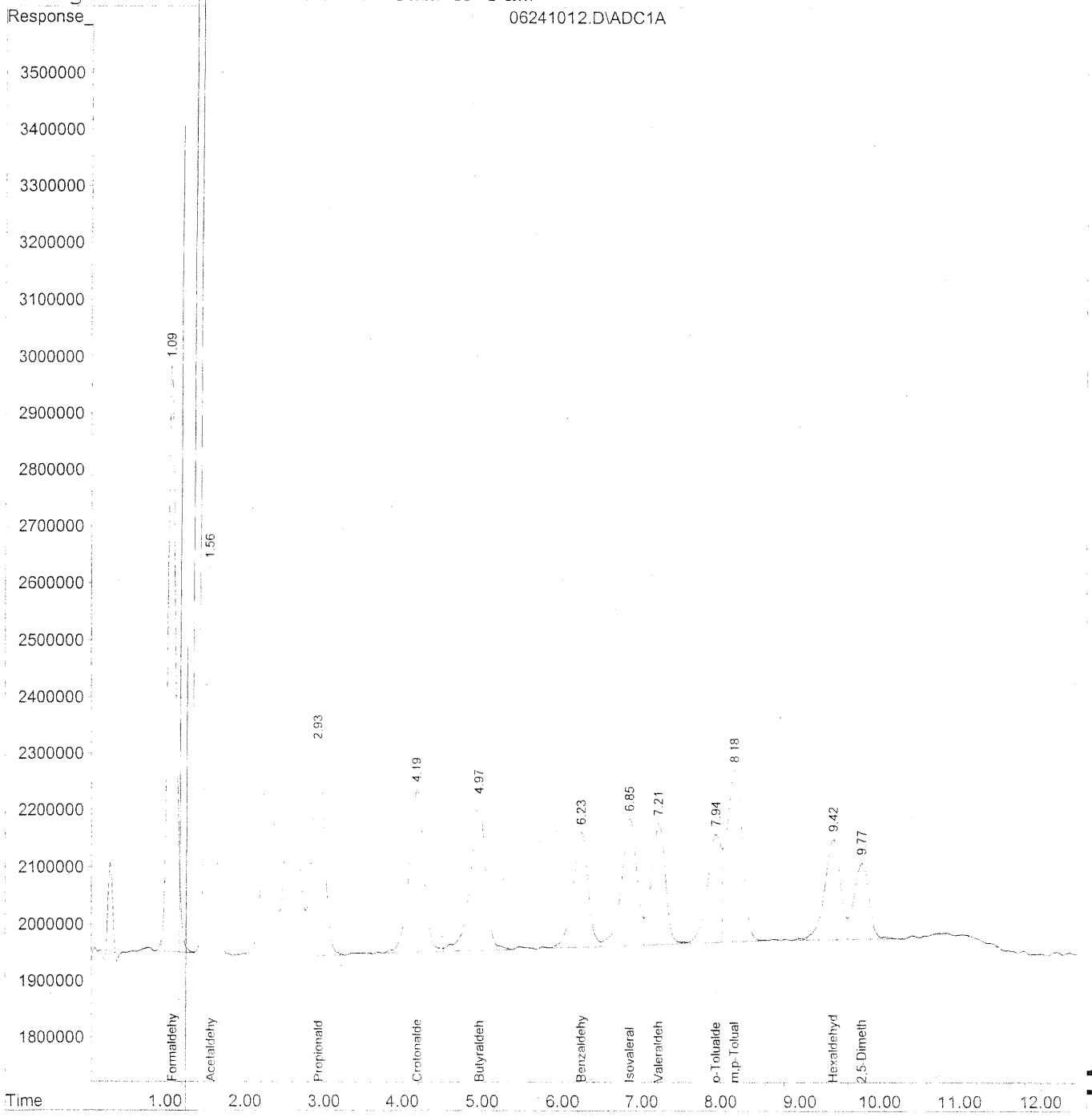
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241012.D Vial: 10
Acq On : 24 Jun 2010 13:52 Operator: MD
Sample : 500ng/ml TO-11A S21-03091008 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 14:27 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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\2010_06\24\06241012.D Vial: 10
Acq On : 24 Jun 2010 13:52 Operator: MD
Sample : 500ng/ml TO-11A S21-03091008 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 14:27 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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	1.09	69824657	508.989	ng/ml
2)	Acetaldehyde	1.57	53171079	489.746	ng/ml
3)	Propionaldehyde	2.93	40786533	513.457	ng/ml
4)	Crotonaldehyde	4.19	36622016	478.446	ng/ml
5)	Butyraldehyde	4.96	35535481	484.599	ng/ml
6)	Benzaldehyde	6.23	25798064	462.892	ng/ml
7)	Isovaleraldehyde	6.85	30475379	479.401	ng/ml
8)	Valeraldehyde	7.21	26332477	456.329	ng/ml
9)	o-Tolualdehyde	7.94	23290726	484.138	ng/ml
10)	m,p-Tolualdehyde	8.18	41386366	982.853	ng/ml
11)	Hexaldehyde	9.42	24685885	465.018	ng/ml
12)	2,5-Dimethylbenzaldehyde	9.78	17286981	435.287	ng/ml

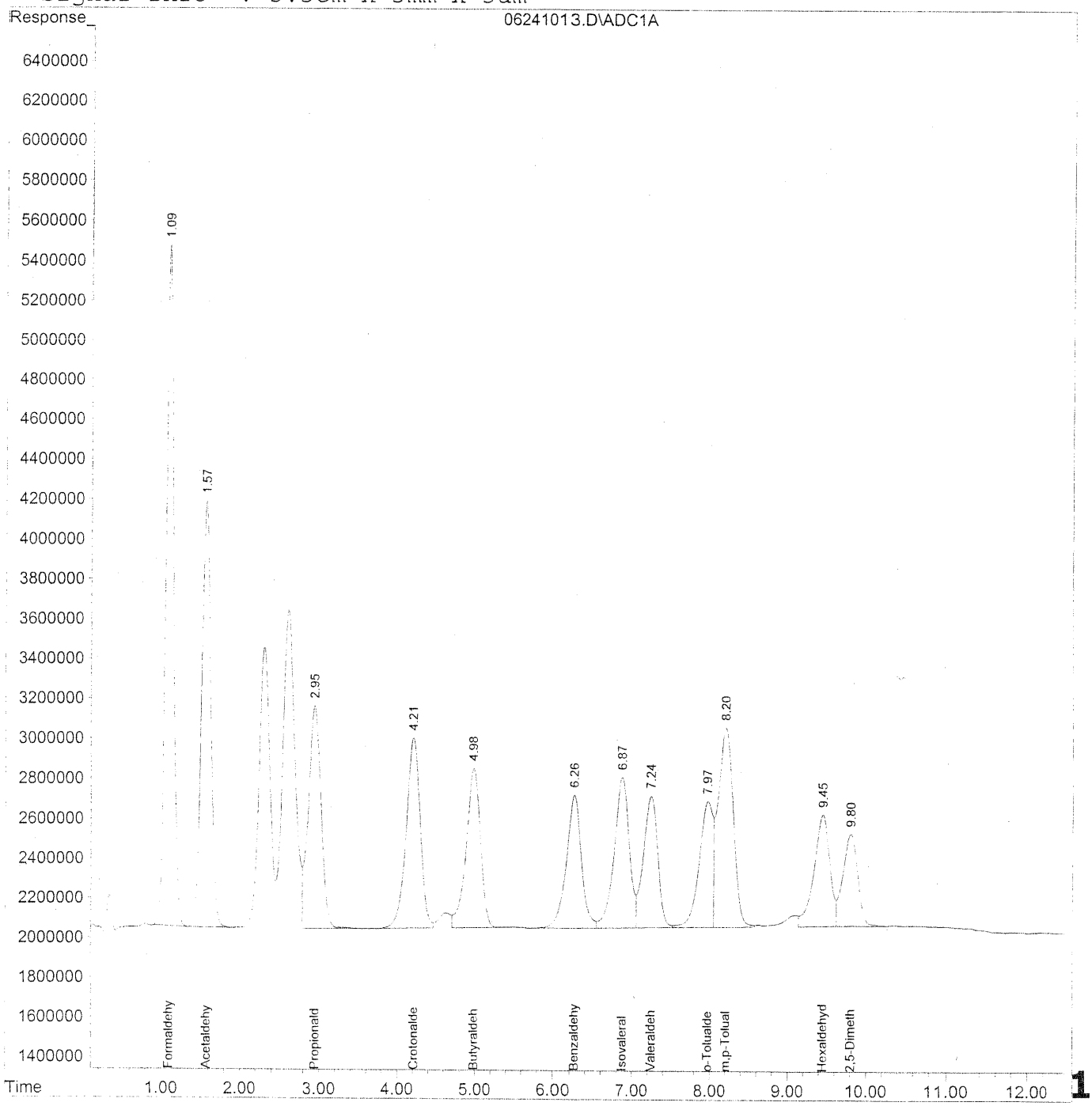
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241013.D Vial: 11
Acq On : 24 Jun 2010 14:06 Operator: MD
Sample : 1500ng/ml TO-11A S21-04211003 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 14:28 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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\2010_06\24\06241013.D Vial: 11
Acq On : 24 Jun 2010 14:06 Operator: MD
Sample : 1500ng/ml TO-11A S21-04211003 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 14:28 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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	1.09	230593016	1683.790	ng/ml
2)	Acetaldehyde	1.57	175356392	1618.937	ng/ml
3)	Propionaldehyde	2.95	125879750	1584.911	ng/ml
4)	Crotonaldehyde	4.21	120696891	1584.647	ng/ml
5)	Butyraldehyde	4.99	103399084	1410.674	ng/ml
6)	Benzaldehyde	6.26	89107201	1610.229	ng/ml
7)	Isovaleraldehyde	6.87	102723576	1624.188	ng/ml
8)	Valeraldehyde	7.24	84987047	1484.702	ng/ml
9)	o-Tolualdehyde	7.97	77943499	1628.761	ng/ml
10)	m,p-Tolualdehyde	8.20	138317069	3296.169	ng/ml
11)	Hexaldehyde	9.44	79547312	1508.911	ng/ml
12)	2,5-Dimethylbenzaldehyde	9.80	60933239	1548.712	ng/ml

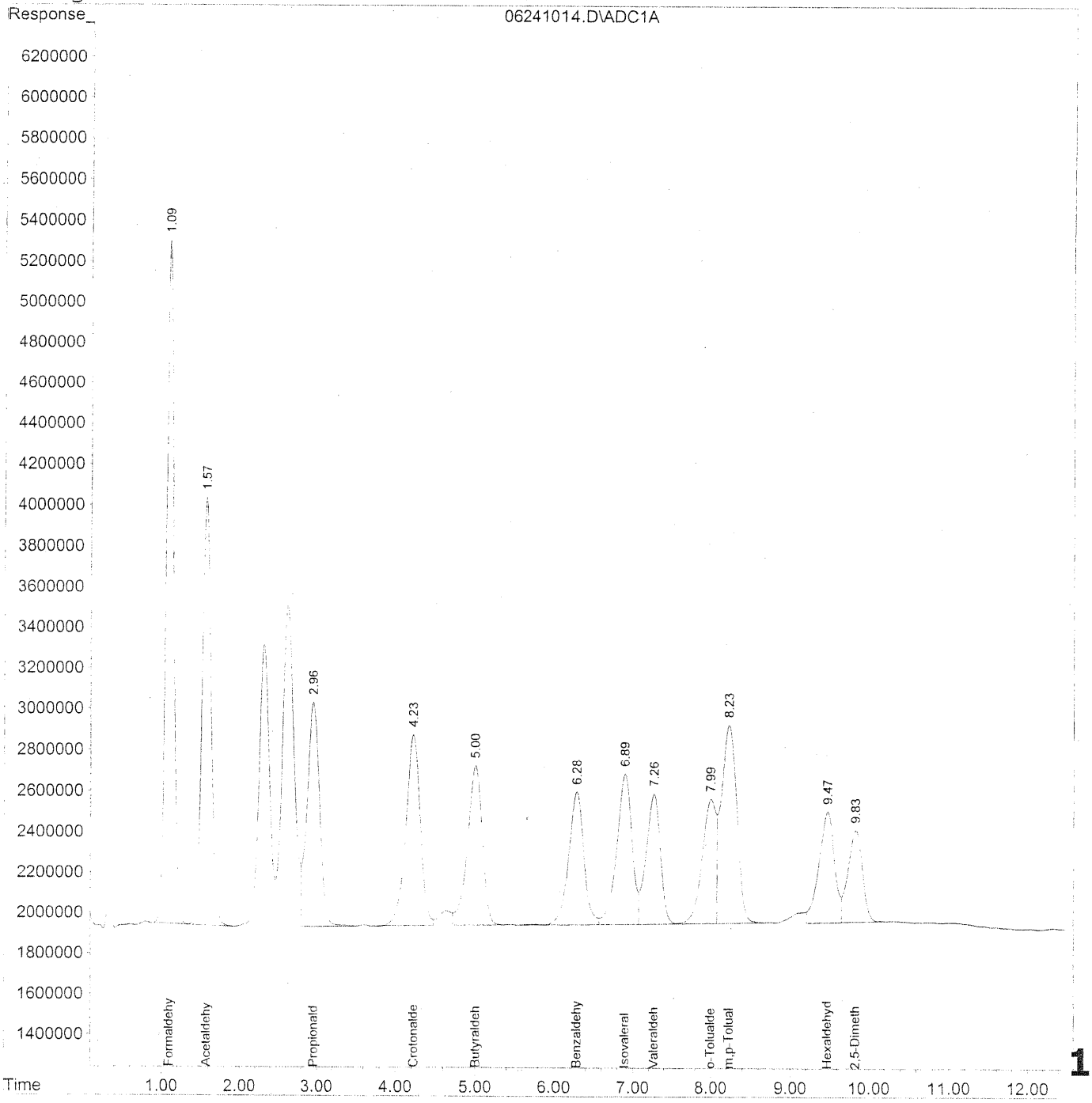
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241014.D Vial: 12
Acq On : 24 Jun 2010 14:19 Operator: MD
Sample : 1500ng/ml TO-11A S21-04211003 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 14:35 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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\2010_06\24\06241014.D Vial: 12
Acq On : 24 Jun 2010 14:19 Operator: MD
Sample : 1500ng/ml TO-11A S21-04211003 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 14:35 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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

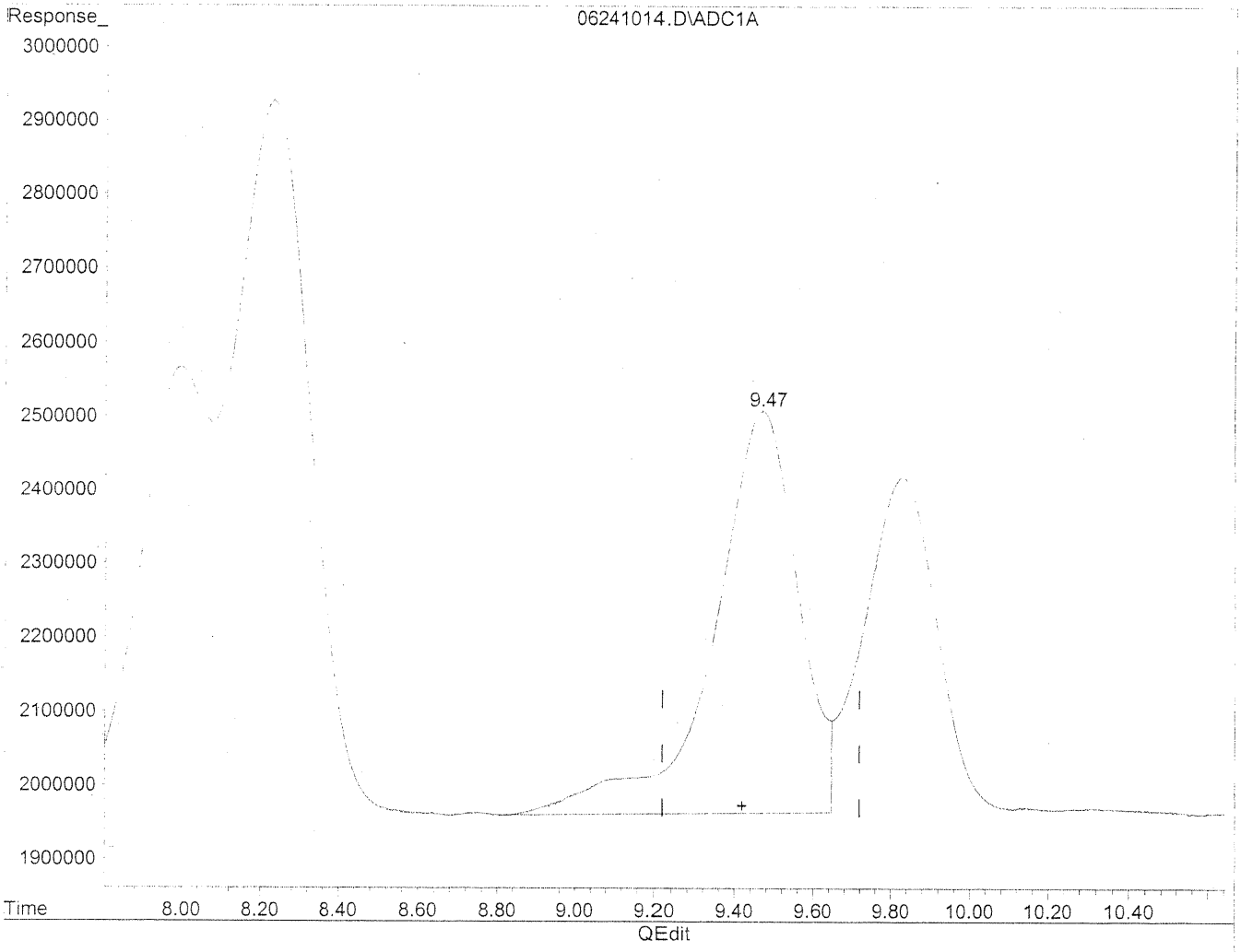
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.09	225421129	1597.103	ng/ml
2)	Acetaldehyde	1.57	172419255	1560.880	ng/ml
3)	Propionaldehyde	2.96	124283598	1542.979	ng/ml
4)	Crotonaldehyde	4.23	119167899	1542.807	ng/ml
5)	Butyraldehyde	5.00	101934633	1411.711	ng/ml
6)	Benzaldehyde	6.28	86843740	1541.016	ng/ml
7)	Isovaleraldehyde	6.89	100332819	1554.218	ng/ml
8)	Valeraldehyde	7.26	81906561	1434.545	ng/ml
9)	o-Tolualdehyde	8.00	73725386	1508.249	ng/ml
10)	m,p-Tolualdehyde	8.23	134049332	3117.524	ng/ml
11)	Hexaldehyde	9.47	75378424	1427.712	ng/mlm
12)	2,5-Dimethylbenzaldehyde	9.83	58178943	1466.799	ng/ml

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241014.D Vial: 12
Acq On : 24 Jun 2010 14:19 Operator: MD
Sample : 1500ng/ml TO-11A S21-04211003 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 14:34 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
Response via : Multiple Level Calibration



(11) Hexaldehyde

9.47min 1538.904ng/ml

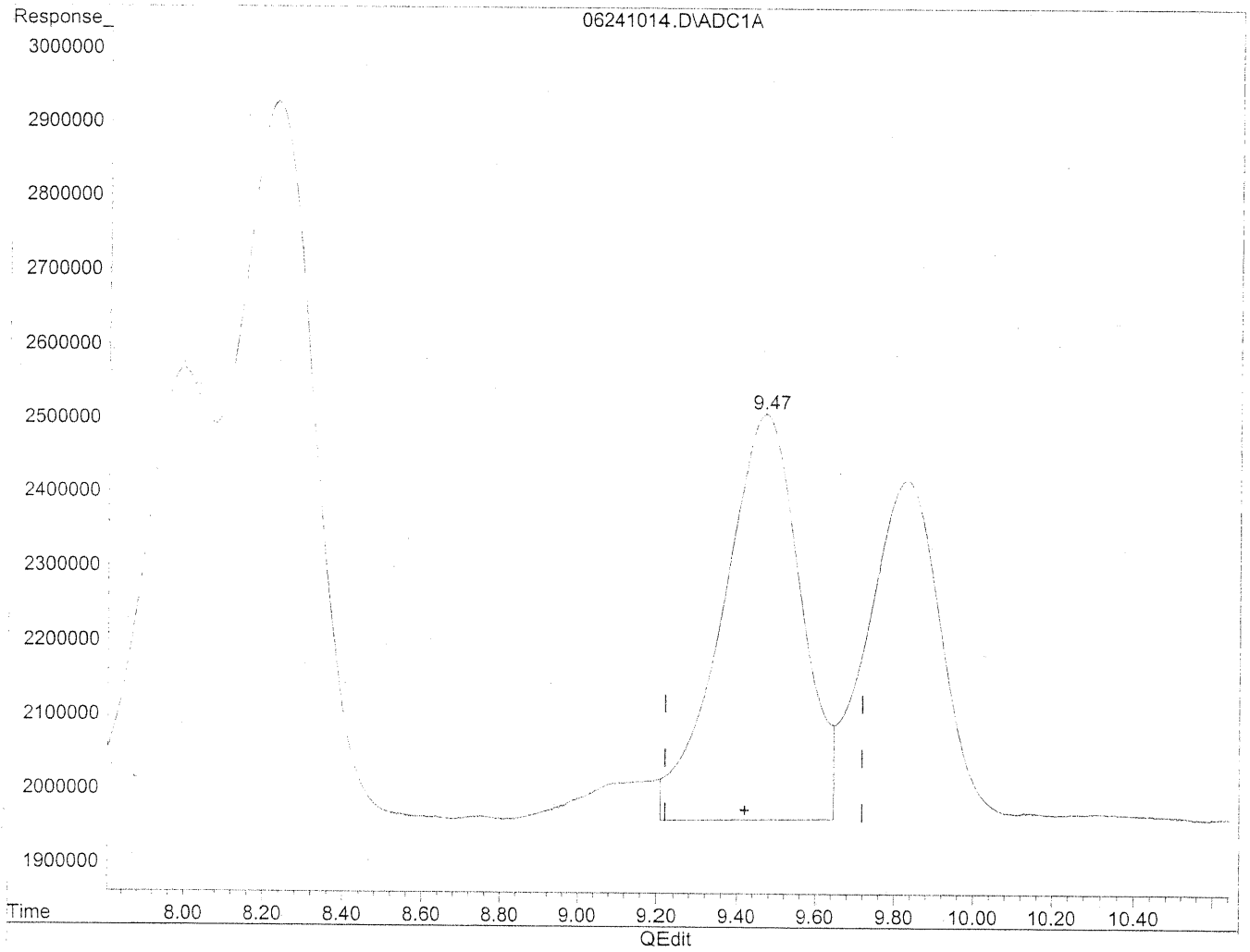
response 81248952

(+) = Expected Retention Time
06241014.D TO110610.M Thu Jun 24 14:34:53 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241014.D Vial: 12
Acq On : 24 Jun 2010 14:19 Operator: MD
Sample : 1500ng/ml TO-11A S21-04211003 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 14:34 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
Response via : Multiple Level Calibration



(11) Hexaldehyde

9.47min 1427.712ng/ml m

response 75378424

6/25/10 6/25/10

189

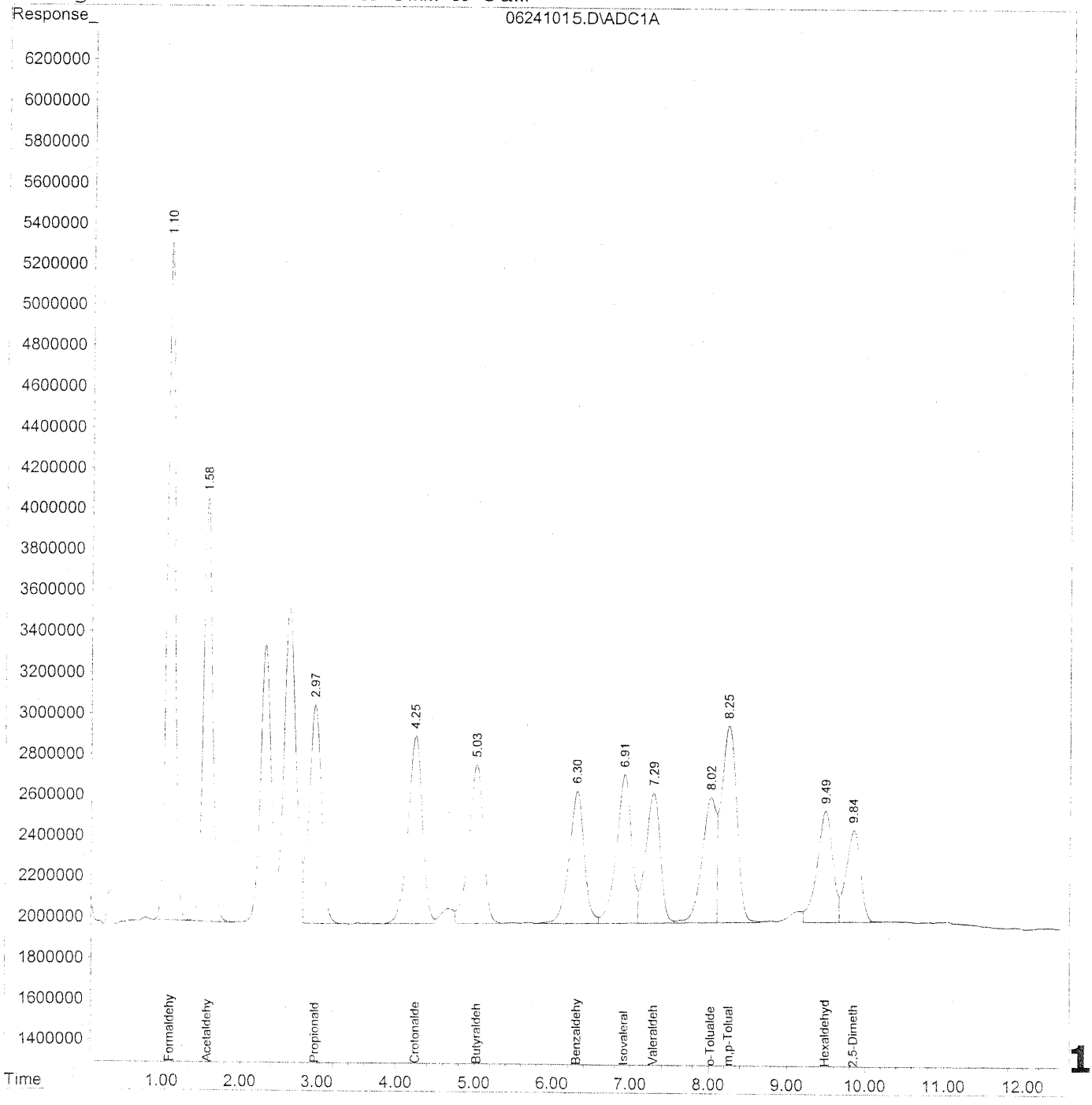
(+) = Expected Retention Time
06241014.D TO110610.M Thu Jun 24 14:35:00 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241015.D Vial: 13
Acq On : 24 Jun 2010 14:33 Operator: MD
Sample : 1500ng/ml TO-11A S21-04211003 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:18 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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\2010_06\24\06241015.D Vial: 13
Acq On : 24 Jun 2010 14:33 Operator: MD
Sample : 1500ng/ml TO-11A S21-04211003 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:18 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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

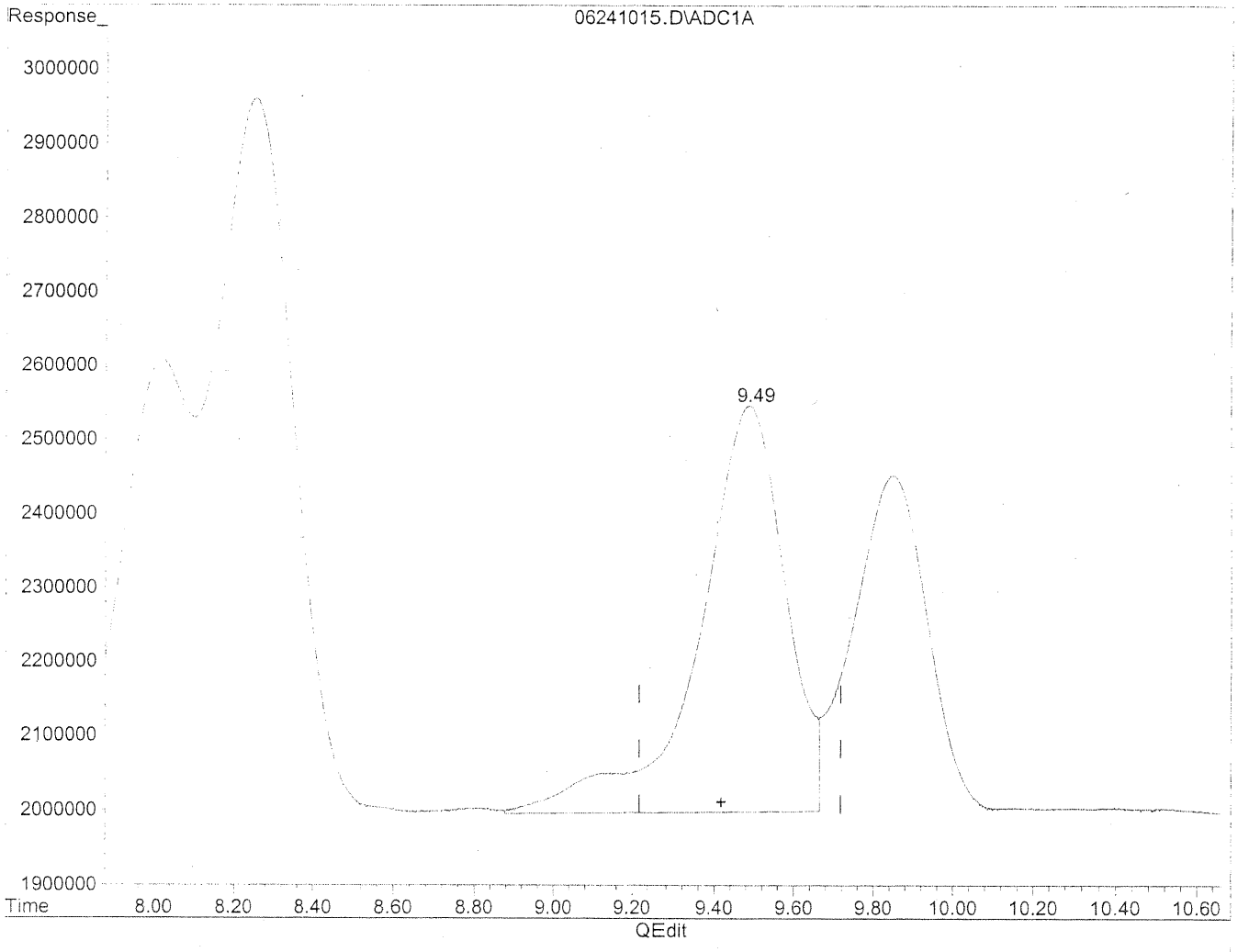
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.10	225129784	1599.924	ng/ml
2)	Acetaldehyde	1.58	170977797	1551.268	ng/ml
3)	Propionaldehyde	2.97	121684007	1513.203	ng/ml
4)	Crotonaldehyde	4.25	116879696	1515.683	ng/ml
5)	Butyraldehyde	5.03	101737279	1411.364	ng/ml
6)	Benzaldehyde	6.31	86160446	1534.026	ng/ml
7)	Isovaleraldehyde	6.92	99712150	1549.385	ng/ml
8)	Valeraldehyde	7.29	81955851	1441.891	ng/ml
9)	o-Tolualdehyde	8.02	76844694	1583.450	ng/ml
10)	m,p-Tolualdehyde	8.25	133062119	3107.415	ng/ml
11)	Hexaldehyde	9.49	76355051	1455.790	ng/mlm
12)	2,5-Dimethylbenzaldehyde	9.85	58479212	1482.951	ng/ml

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241015.D Vial: 13
Acq On : 24 Jun 2010 14:33 Operator: MD
Sample : 1500ng/ml TO-11A S21-04211003 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:17 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
Response via : Multiple Level Calibration



(11) Hexaldehyde

9.49min 1575.238ng/ml

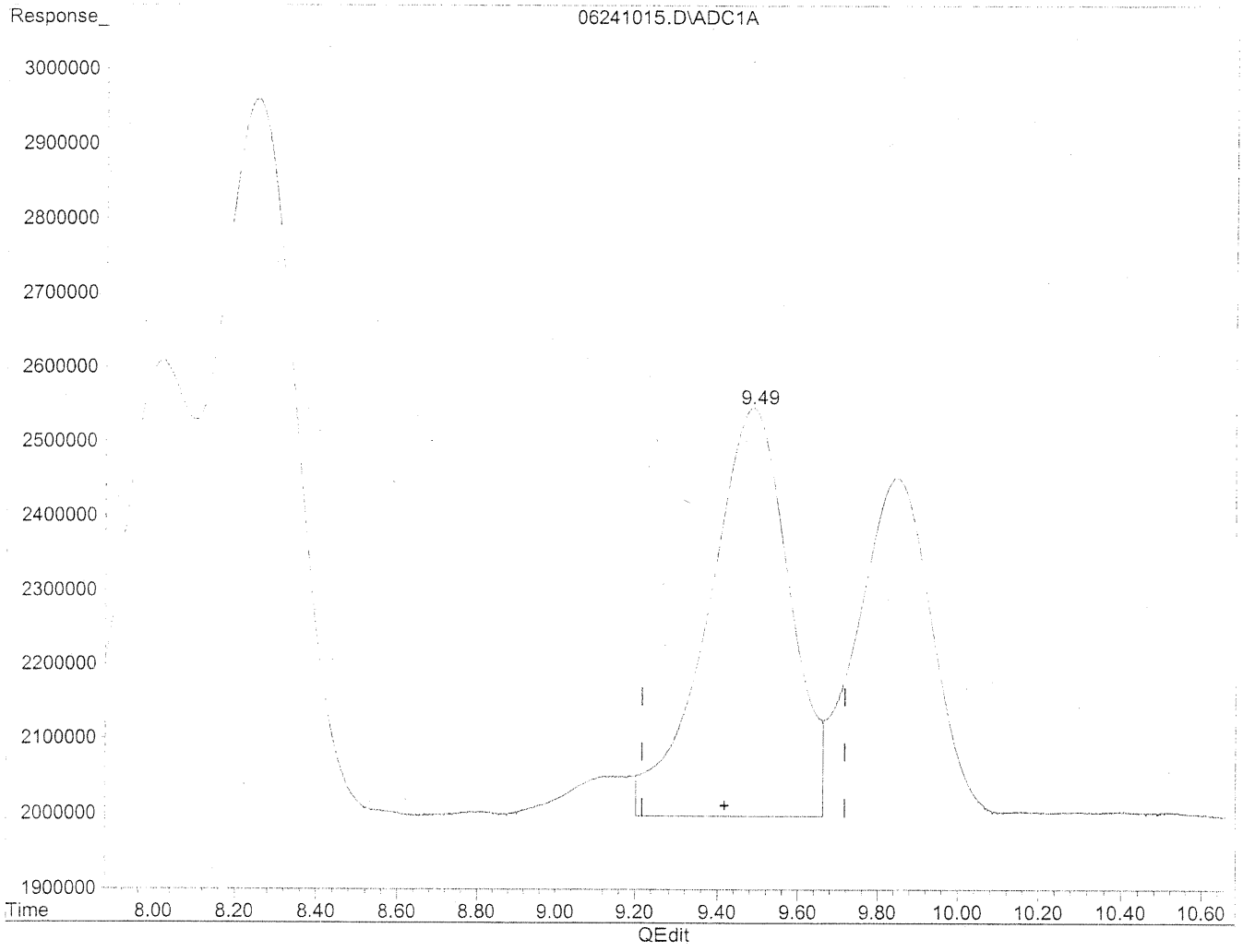
response 82620034

(+) = Expected Retention Time
06241015.D TO110610.M Thu Jun 24 15:18:09 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241015.D Vial: 13
Acq On : 24 Jun 2010 14:33 Operator: MD
Sample : 1500ng/ml TO-11A S21-04211003 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:17 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
Response via : Multiple Level Calibration



(11) Hexaldehyde

9.49min 1455.790ng/ml m

response 76355051

4/28/10 6/25/10

193

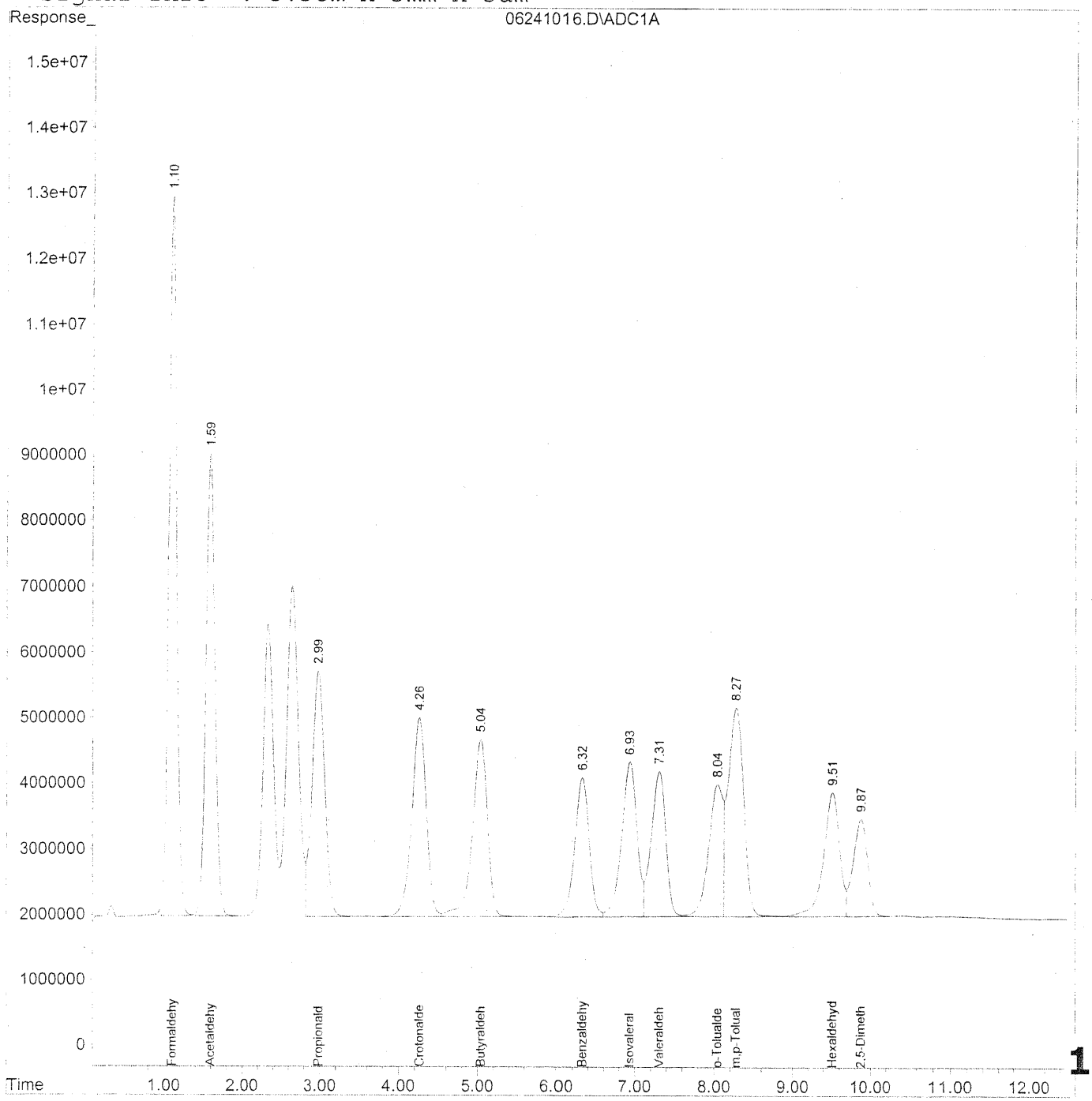
(+) = Expected Retention Time
06241015.D TO110610.M Thu Jun 24 15:18:19 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241016.D Vial: 14
Acq On : 24 Jun 2010 14:46 Operator: MD
Sample : 5000ng/ml TO-11A S21-03091001 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:18 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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



194

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2010_06\24\06241016.D Vial: 14
Acq On : 24 Jun 2010 14:46 Operator: MD
Sample : 5000ng/ml TO-11A S21-03091001 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:18 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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	1.10	752167389	5351.489	ng/ml
2)	Acetaldehyde	1.59	565379435	5137.177	ng/ml
3)	Propionaldehyde	2.99	421481275	5253.669	ng/ml
4)	Crotonaldehyde	4.26	387209880	5032.362	ng/ml
5)	Butyraldehyde	5.04	360334815	5002.375	ng/ml
6)	Benzaldehyde	6.33	271843062	4848.677	ng/ml
7)	Isovaleraldehyde	6.94	318898458	4963.009	ng/ml
8)	Valeraldehyde	7.31	283824088	5000.748	ng/ml
9)	o-Tolualdehyde	8.04	245490595	5052.697	ng/ml
10)	m,p-Tolualdehyde	8.28	439534493	10285.324	ng/ml
11)	Hexaldehyde	9.51	269904263	5152.055	ng/ml
12)	2,5-Dimethylbenzaldehyde	9.87	191751893	4869.948	ng/ml

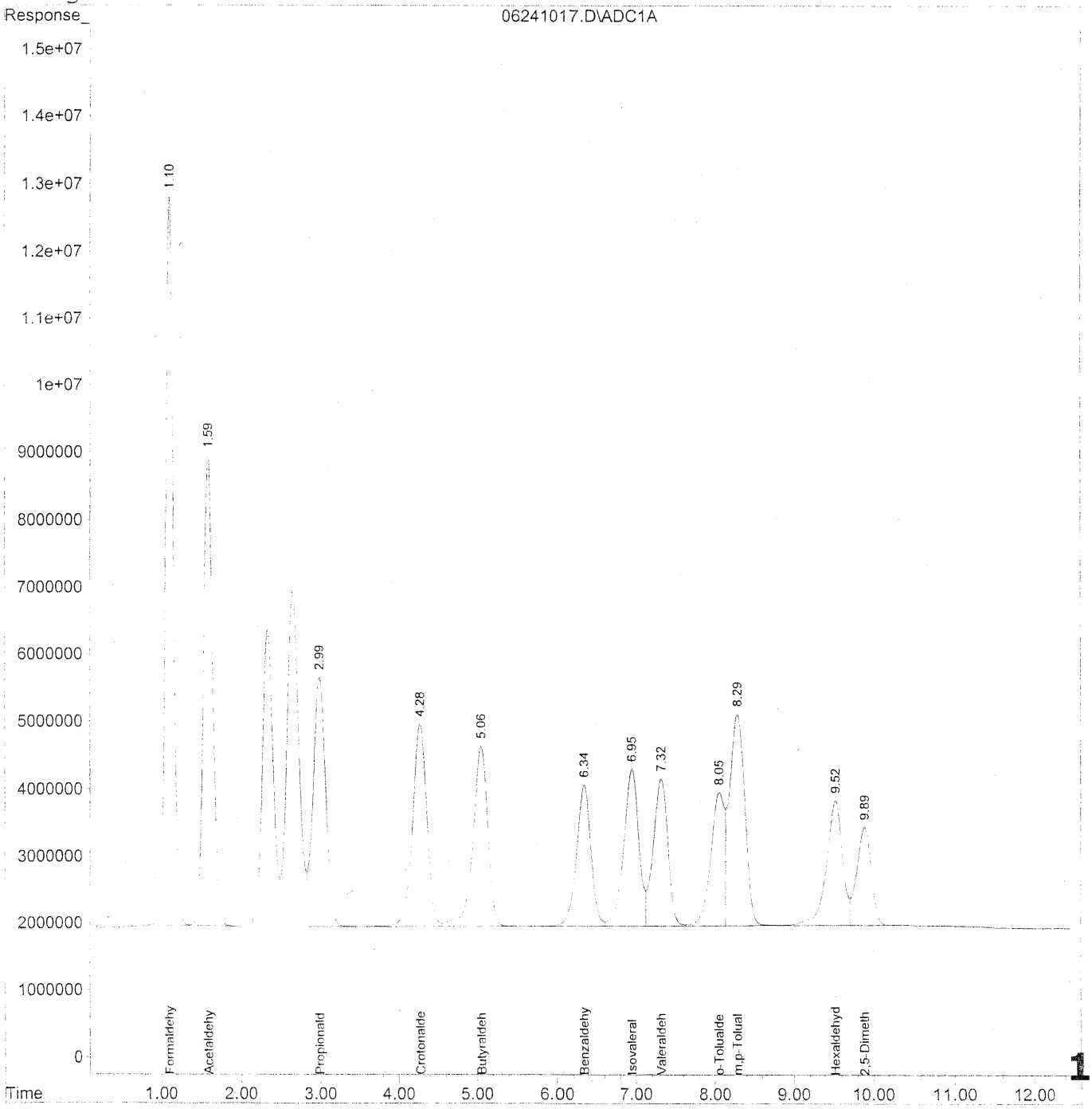
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241017.D Vial: 15
Acq On : 24 Jun 2010 15:00 Operator: MD
Sample : 5000ng/ml TO-11A S21-03091001 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:19 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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\2010_06\24\06241017.D Vial: 15
Acq On : 24 Jun 2010 15:00 Operator: MD
Sample : 5000ng/ml TO-11A S21-03091001 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:19 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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	1.10	741687155	5203.762	ng/ml
2)	Acetaldehyde	1.59	560621731	5066.149	ng/ml
3)	Propionaldehyde	2.99	418074248	5158.856	ng/ml
4)	Crotonaldehyde	4.28	384212117	4986.946	ng/ml
5)	Butyraldehyde	5.06	359182440	4985.904	ng/ml
6)	Benzaldehyde	6.34	271646285	4874.674	ng/ml
7)	Isovaleraldehyde	6.95	316239705	4928.924	ng/ml
8)	Valeraldehyde	7.32	281768408	4964.380	ng/ml
9)	o-Tolualdehyde	8.06	241706915	4964.357	ng/ml
10)	m,p-Tolualdehyde	8.29	435865650	10141.599	ng/ml
11)	Hexaldehyde	9.52	266567105	5057.593	ng/ml
12)	2,5-Dimethylbenzaldehyde	9.89	189585119	4840.097	ng/ml

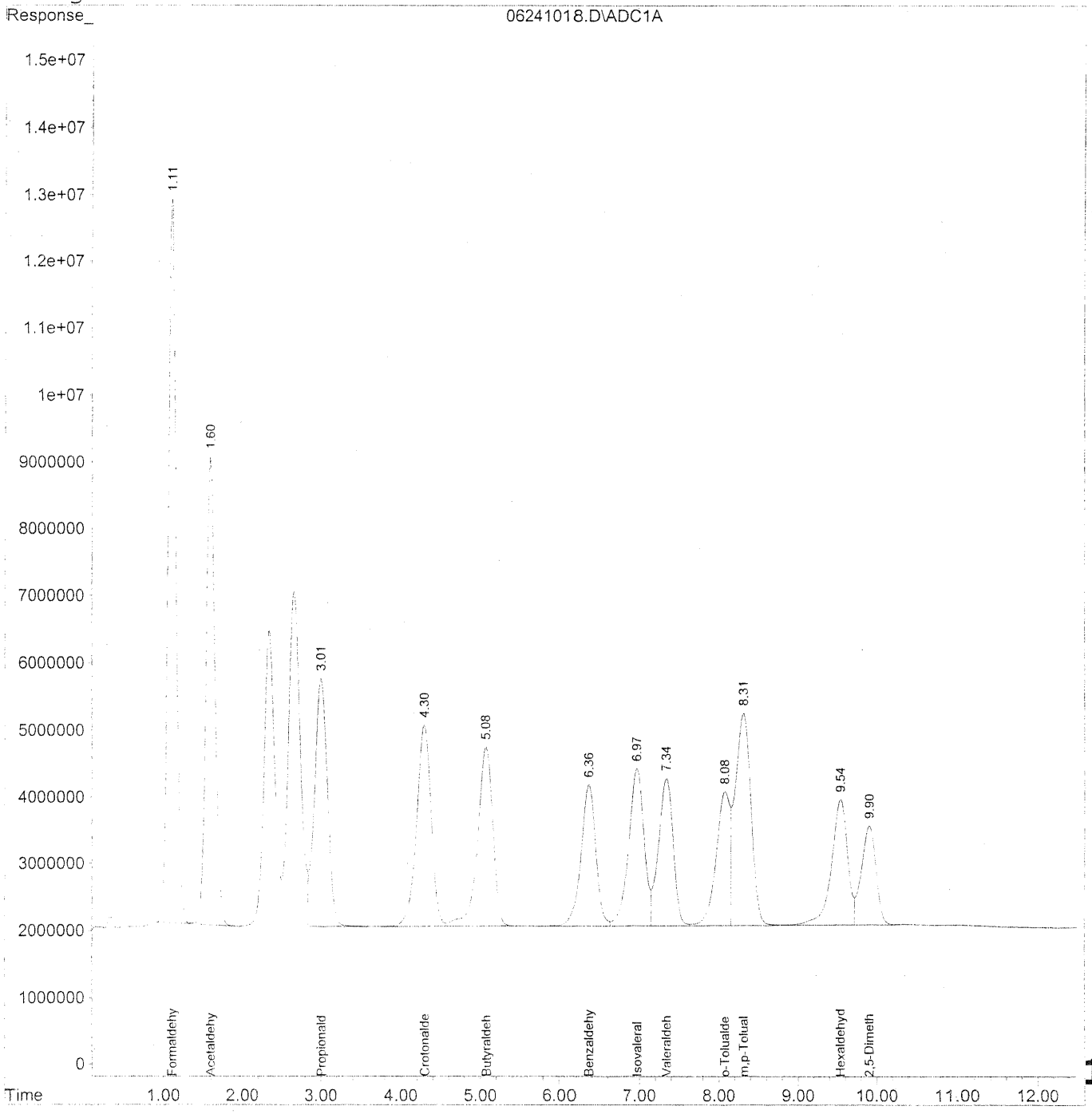
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241018.D Vial: 16
Acq On : 24 Jun 2010 15:13 Operator: MD
Sample : 5000ng/ml TO-11A S21-03091001 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:29 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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\2010_06\24\06241018.D Vial: 16
Acq On : 24 Jun 2010 15:13 Operator: MD
Sample : 5000ng/ml TO-11A S21-03091001 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:29 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
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

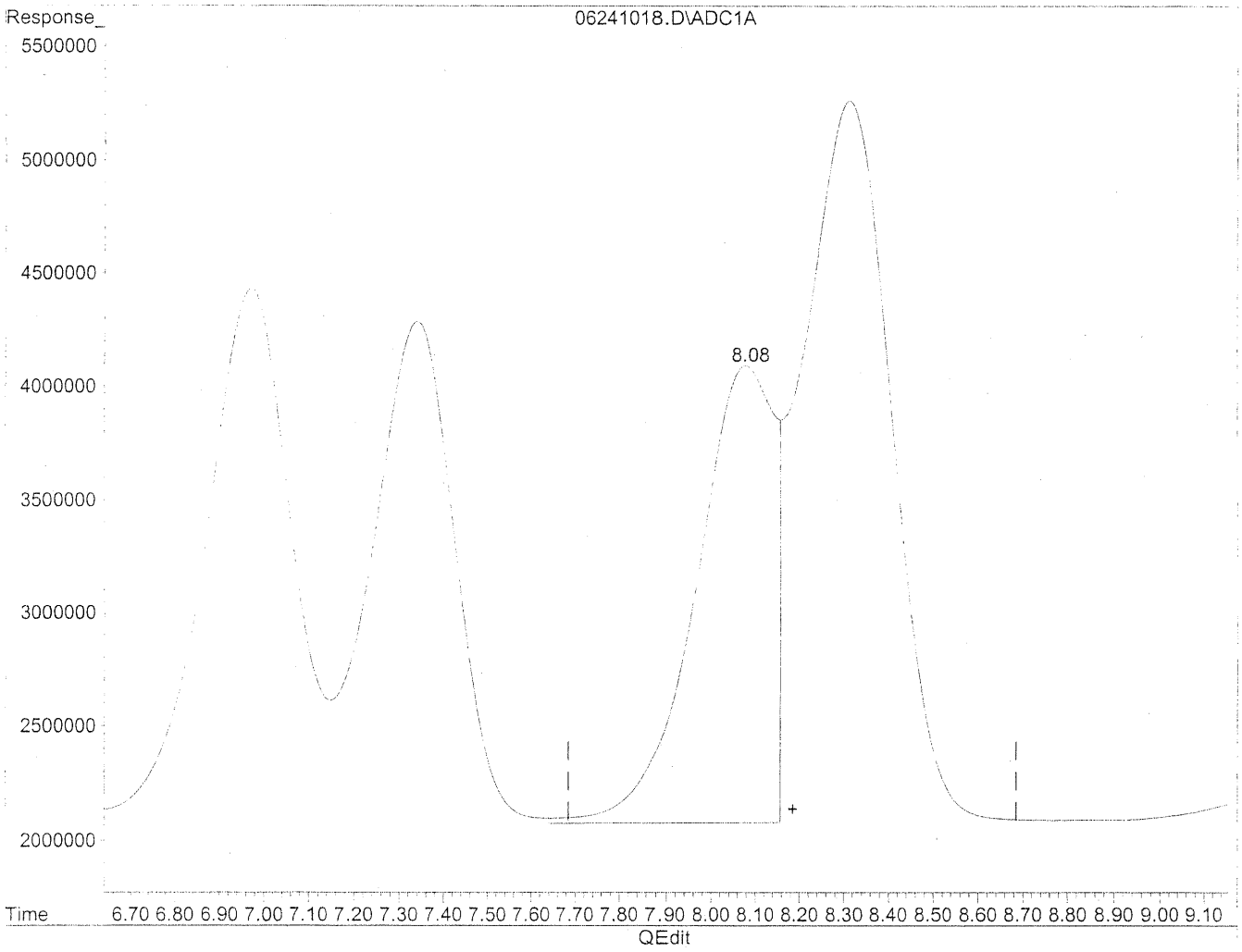
	Compound	R.T.	Response	Conc	Units
	Target Compounds				
1)	Formaldehyde	1.11	735352713	5166.917	ng/ml
2)	Acetaldehyde	1.60	560284478	5067.458	ng/ml
3)	Propionaldehyde	3.01	420570615	5194.027	ng/ml
4)	Crotonaldehyde	4.30	385901475	5012.774	ng/ml
5)	Butyraldehyde	5.08	359960551	4998.304	ng/ml
6)	Benzaldehyde	6.36	273410909	4906.686	ng/ml
7)	Isovaleraldehyde	6.97	319488895	4983.696	ng/ml
8)	Valeraldehyde	7.34	284129216	5009.603	ng/ml
9)	o-Tolualdehyde	8.08	244039778	5020.073	ng/ml
10)	m,p-Tolualdehyde	8.31	439548558	10236.030	ng/mlm
11)	Hexaldehyde	9.54	269656485	5122.695	ng/ml
12)	2,5-Dimethylbenzaldehyde	9.90	191978970	4906.640	ng/ml

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

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241018.D Vial: 16
Acq On : 24 Jun 2010 15:13 Operator: MD
Sample : 5000ng/ml TO-11A S21-03091001 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:29 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

8.08min 5683.100ng/ml

response 244039778

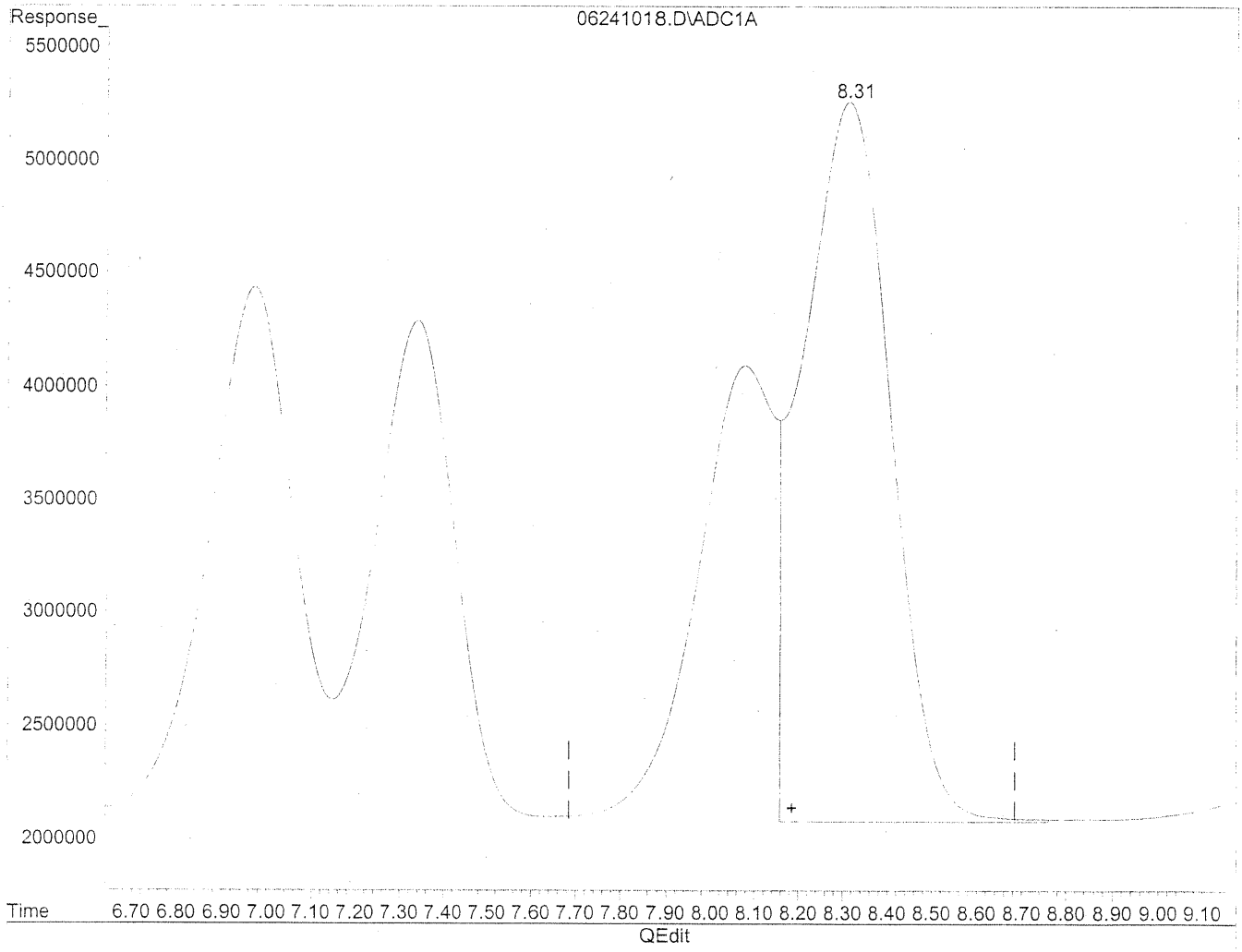
200

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241018.D Vial: 16
Acq On : 24 Jun 2010 15:13 Operator: MD
Sample : 5000ng/ml TO-11A S21-03091001 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:29 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 13:47:22 2010
Response via : Multiple Level Calibration



(10) m,p-Tolualdehyde

8.31min 10236.030ng/ml m

response 439548558

6/25/10 6/25/10

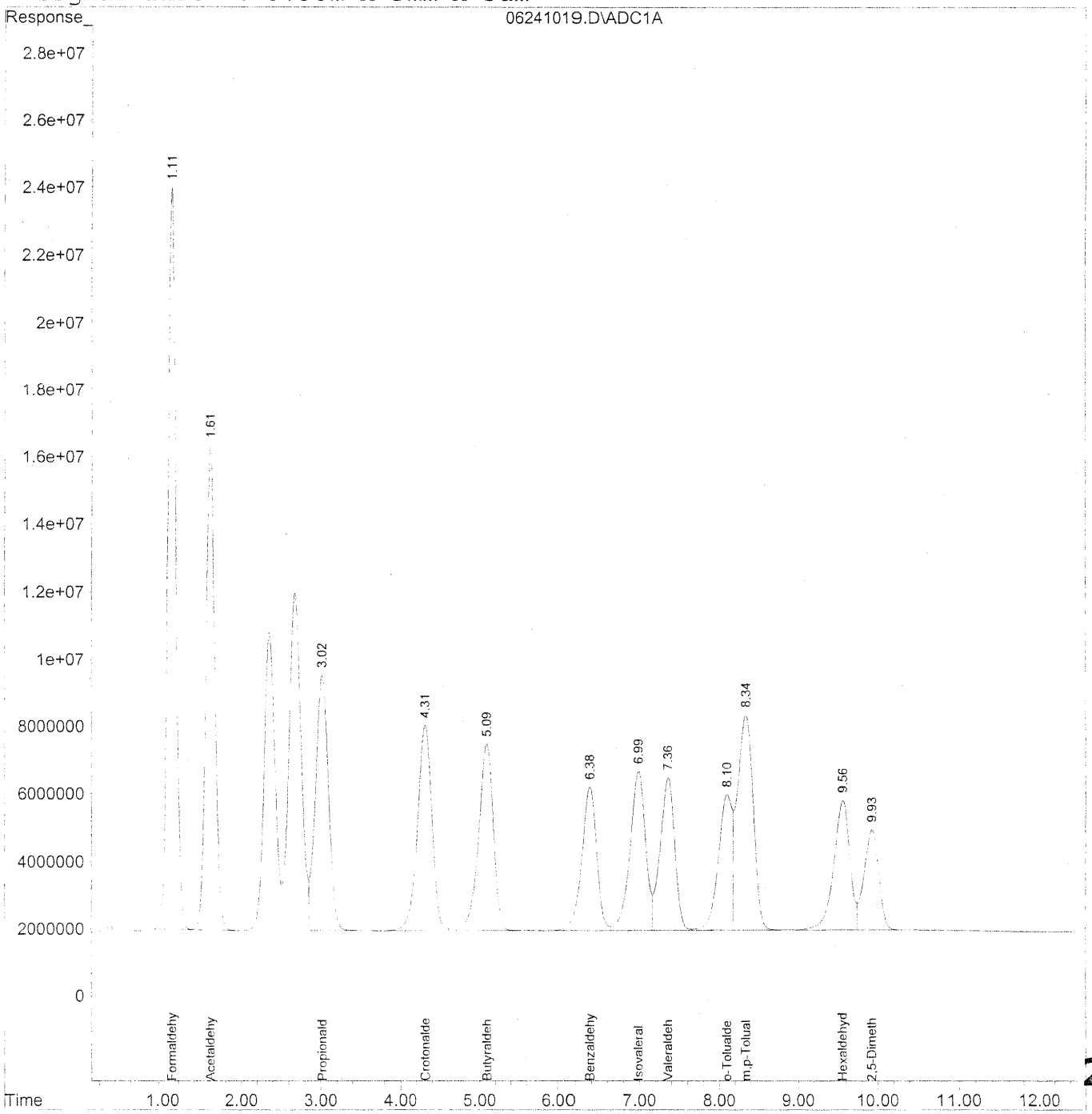
(+) = Expected Retention Time
06241018.D TO110610.M Thu Jun 24 15:29:31 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241019.D Vial: 17
Acq On : 24 Jun 2010 15:27 Operator: MD
Sample : 10000ng/ml TO-11A S21-03091007 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:49 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
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



202

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2010_06\24\06241019.D Vial: 17
Acq On : 24 Jun 2010 15:27 Operator: MD
Sample : 10000ng/ml TO-11A S21-03091007 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:49 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
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	1.11	1506388672	10596.051	ng/ml
2)	Acetaldehyde	1.61	1143475045	10345.478	ng/ml
3)	Propionaldehyde	3.02	857670773	10590.811	ng/ml
4)	Crotonaldehyde	4.31	774502971	10060.289	ng/ml
5)	Butyraldehyde	5.09	723112269	10040.544	ng/ml
6)	Benzaldehyde	6.38	541585412	9715.525	ng/ml
7)	Isovaleraldehyde	6.99	637094049	9934.041	ng/ml
8)	Valeraldehyde	7.36	579085929	10206.912	ng/ml
9)	o-Tolualdehyde	8.10	490639974	10091.596	ng/ml
10)	m,p-Tolualdehyde	8.33	882689956	20549.830	ng/ml
11)	Hexaldehyde	9.56	537444487	10206.220	ng/ml
12)	2,5-Dimethylbenzaldehyde	9.93	383332870	9792.931	ng/ml

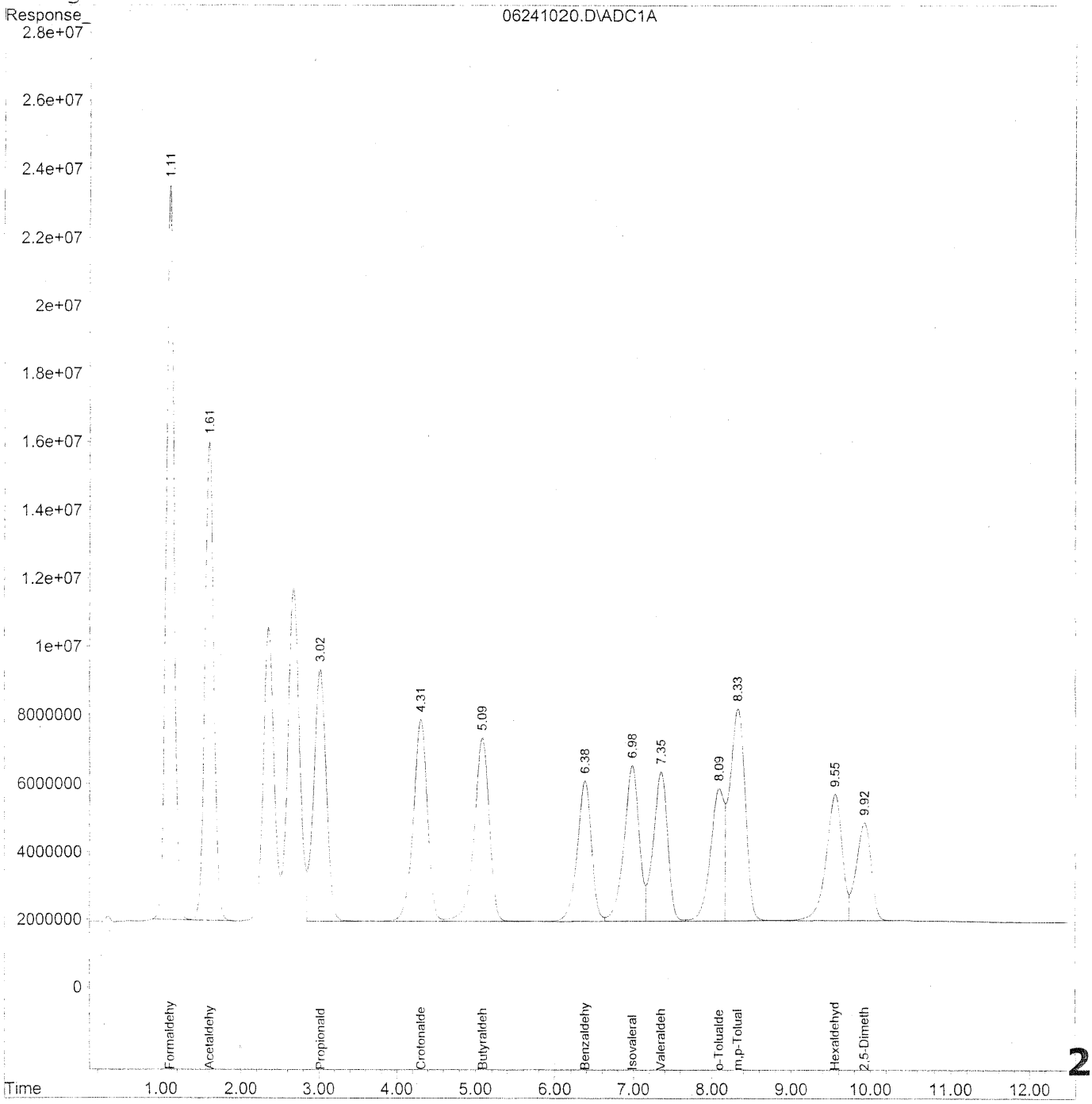
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241020.D Vial: 18
Acq On : 24 Jun 2010 15:40 Operator: MD
Sample : 10000ng/ml TO-11A S21-03091007 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:56 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
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



204

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2010_06\24\06241020.D Vial: 18
Acq On : 24 Jun 2010 15:40 Operator: MD
Sample : 10000ng/ml TO-11A S21-03091007 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 15:56 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
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	1.11	1465408979	10206.404	ng/ml
2)	Acetaldehyde	1.61	1108342081	9970.208	ng/ml
3)	Propionaldehyde	3.02	834978009	10210.056	ng/ml
4)	Crotonaldehyde	4.31	755899749	9808.789	ng/ml
5)	Butyraldehyde	5.09	707762167	9820.769	ng/ml
6)	Benzaldehyde	6.38	527522590	9508.333	ng/ml
7)	Isovaleraldehyde	6.98	619285794	9666.989	ng/ml
8)	Valeraldehyde	7.35	564056973	9907.845	ng/ml
9)	o-Tolualdehyde	8.09	475987748	9775.302	ng/ml
10)	m,p-Tolualdehyde	8.33	860464776	19941.040	ng/ml
11)	Hexaldehyde	9.55	525177534	9939.107	ng/ml
12)	2,5-Dimethylbenzaldehyde	9.92	373842408	9583.554	ng/ml

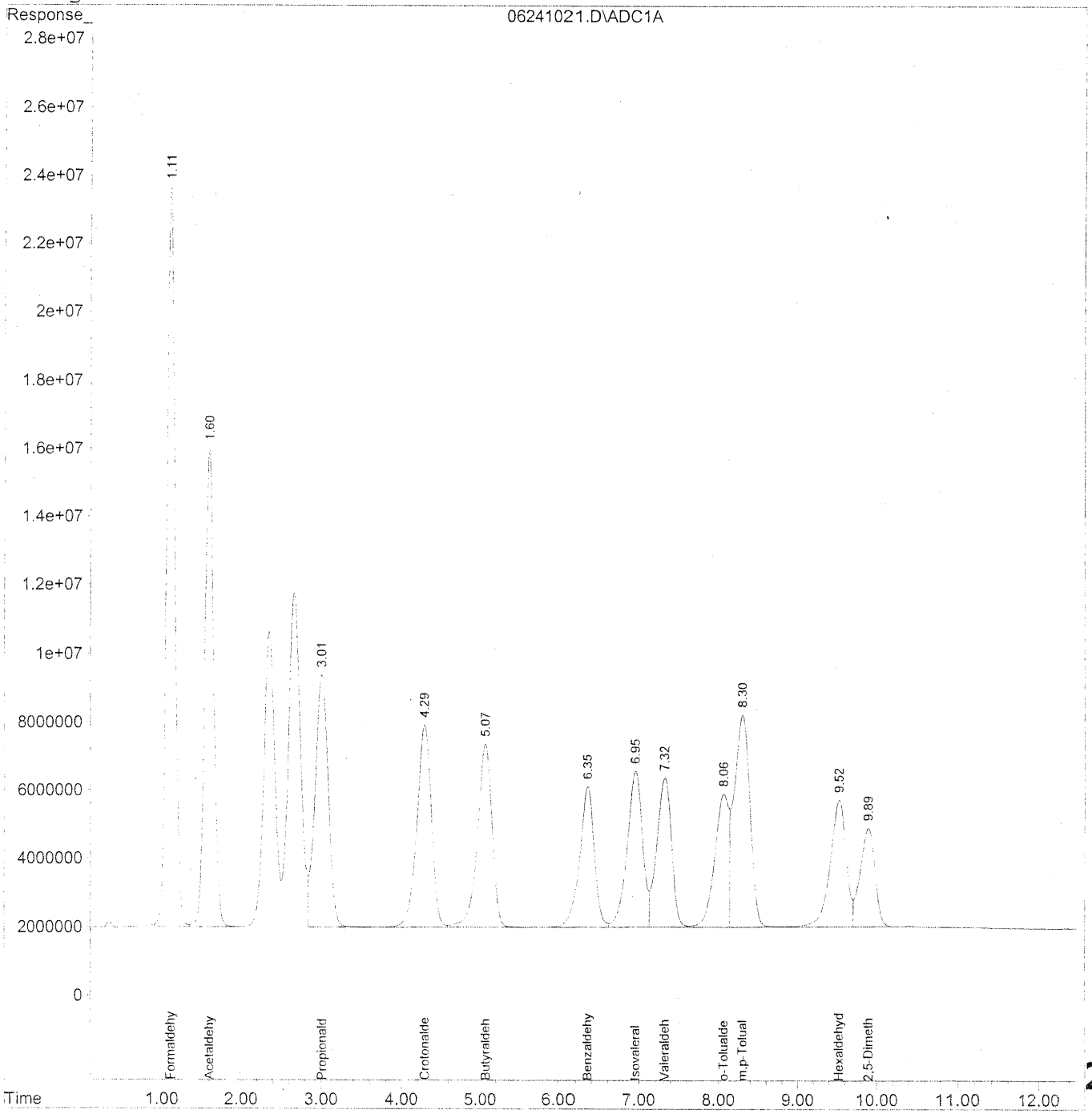
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241021.D Vial: 19
Acq On : 24 Jun 2010 15:54 Operator: MD
Sample : 10000ng/ml TO-11A S21-03091007 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:08 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
Response via : Multiple Level Calibration
DataAcq Meth : TO11S.M

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



206

Quantitation Report (QT Reviewed)

Data File : J:\LC01\DATA\TO11\2010_06\24\06241021.D Vial: 19
Acq On : 24 Jun 2010 15:54 Operator: MD
Sample : 10000ng/ml TO-11A S21-03091007 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:08 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
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	1.11	1476855713	10310.653	ng/ml
2)	Acetaldehyde	1.60	1110424466	10015.318	ng/ml
3)	Propionaldehyde	3.01	832841780	10207.538	ng/ml
4)	Crotonaldehyde	4.29	753218148	9793.694	ng/ml
5)	Butyraldehyde	5.07	704333220	9790.568	ng/ml
6)	Benzaldehyde	6.35	525054783	9483.885	ng/ml
7)	Isovaleraldehyde	6.96	618766885	9681.316	ng/ml
8)	Valeraldehyde	7.32	562282799	9898.457	ng/ml
9)	o-Tolualdehyde	8.06	474574515	9770.780	ng/ml
10)	m,p-Tolualdehyde	8.30	859080340	19951.774	ng/ml
11)	Hexaldehyde	9.52	521619735	9890.910	ng/ml
12)	2,5-Dimethylbenzaldehyde	9.89	372739681	9574.697	ng/ml

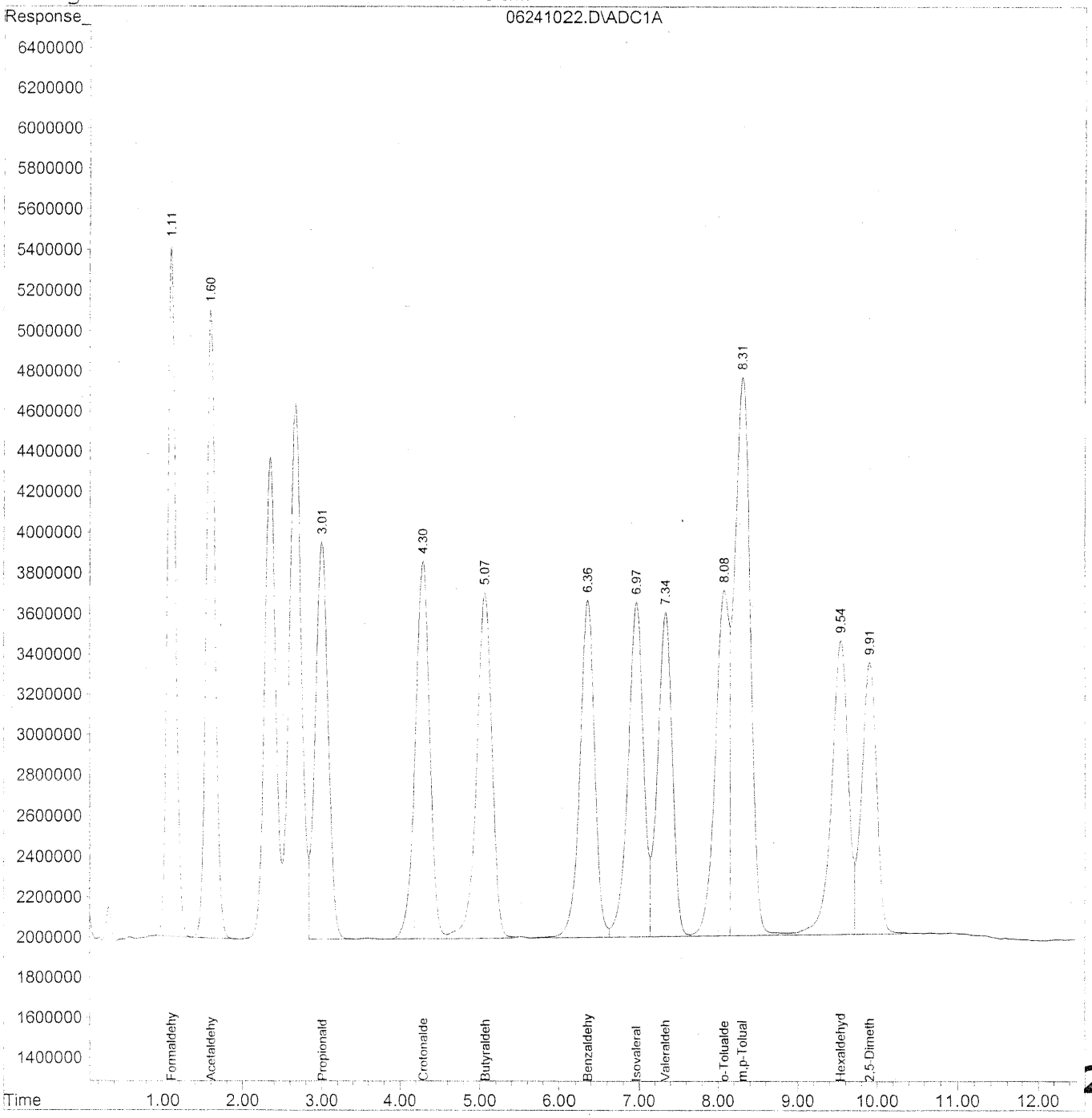
(f)=RT Delta > 1/2 Window (m)=manual int.

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\24\06241022.D Vial: 20
Acq On : 24 Jun 2010 16:07 Operator: MD
Sample : ~1500ng/ml TO-11A ICV S21-03091003 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:56 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
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\2010_06\24\06241022.D Vial: 20
Acq On : 24 Jun 2010 16:07 Operator: MD
Sample : ~1500ng/ml TO-11A ICV S21-03091003 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 24 16:56 19110 Quant Results File: TO110610.RES

Quant Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Thu Jun 24 15:49:23 2010
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	1.11	234075256	1634.767	ng/ml
2)	Acetaldehyde	1.60	246037651	2231.481	ng/ml
3)	Propionaldehyde	3.01	222241976	2726.362	ng/ml
4)	Crotonaldehyde	4.30	237673277	3109.807	ng/ml
5)	Butyraldehyde	5.08	226379316	3157.922	ng/ml
6)	Benzaldehyde	6.36	214780298	3906.517	ng/ml
7)	Isovaleraldehyde	6.97	225723928	3536.309	ng/ml
8)	Valeraldehyde	7.34	207183915	3636.017	ng/ml
9)	o-Tolualdehyde	8.08	205946824	4264.643	ng/ml
10)	m,p-Tolualdehyde	8.31	387123455	8965.224	ng/ml
11)	Hexaldehyde	9.54	207686169	3950.944	ng/ml
12)	2,5-Dimethylbenzaldehyde	9.91	176117645	4498.740	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: 06/12/11

	MW	Aldehyde-DNPH MW*	Manufacturer Prepared Concentration as Aldehyde-DNPH (ug/mL)	Calculated Concentration as Aldehyde (ug/mL)	ICV S21-03091003 (nominal ng/mL)	ICV S21-03091003 (Actual, ng/mL)	% Diff
Formaldehyde	30.03	210.03	100	14.30	1430	1634.77	14.32%
Acetaldehyde	44.05	224.05	100.2	19.70	1970	2231.48	13.27%
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	2726.36	11.55%
Crotonaldehyde	70.09	250.09	100.2	28.08	2808	3109.81	10.75%
Butyraldehyde	72.11	252.11	100	28.60	2860	3157.92	10.42%
Benzaldehyde	106.12	286.12	100	37.09	3709	3906.52	5.33%
Isovaleraldehyde	86.13	266.13	100.2	32.43	3243	3536.31	9.04%
Valeraldehyde	86.13	266.13	100.1	32.40	3240	3636.02	12.22%
o-Tolualdehyde	120.15	300.15	100.1	40.07	4007	4264.64	6.43%
m,p-Tolualdehyde	120.15	300.15	100.3	80.30	8030	8965.22	11.65%
Hexaldehyde	100.16	280.16	100.3	35.86	3586	3950.94	10.18%
2,5-Dimethylbenzaldehyde	134.18	314.18	100.3	42.84	4284	4498.74	5.01%

(* 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

QC

06/28/10

CCV S21-03091001	5173.60
ACTUAL	5000.00
%DIFFERENCE	3.5%
CCV S21-03091001	5075.04
ACTUAL	5000.00
%DIFFERENCE	1.5%
CCV S21-03091001	5069.14
ACTUAL	5000.00
%DIFFERENCE	1.4%
CCV S21-03091001	5267.29
ACTUAL	5000.00
%DIFFERENCE	5.3%
CCV S21-03091001	5021.16
ACTUAL	5000.00
%DIFFERENCE	0.4%
End std S21-03091001	5129.65
ACTUAL	5000.00
%DIFFERENCE	2.6%

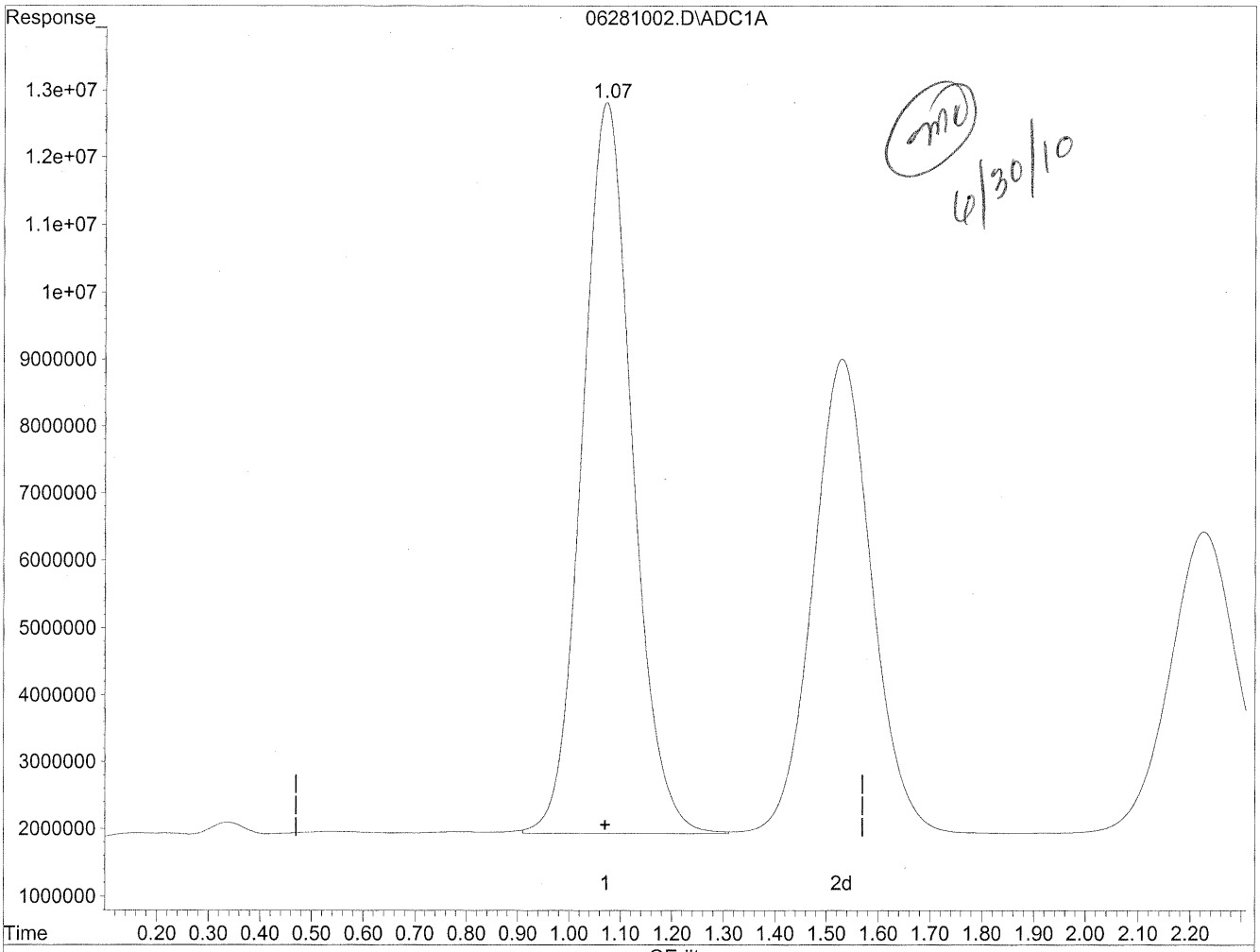
06/29/10

CCV S21-04211003	1620.81
ACTUAL	1500.00
%DIFFERENCE	8.1%
End std S21-04211003	1580.91
ACTUAL	1500.00
%DIFFERENCE	5.4%

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281002.D Vial: 2
Acq On : 28 Jun 2010 10:20 Operator: MD
Sample : 5000ng/ml TO-11A S21-030910001 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 12:58 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.07min 5173.597ng/ml

response 740784986

6/30/10

6/30/10

213

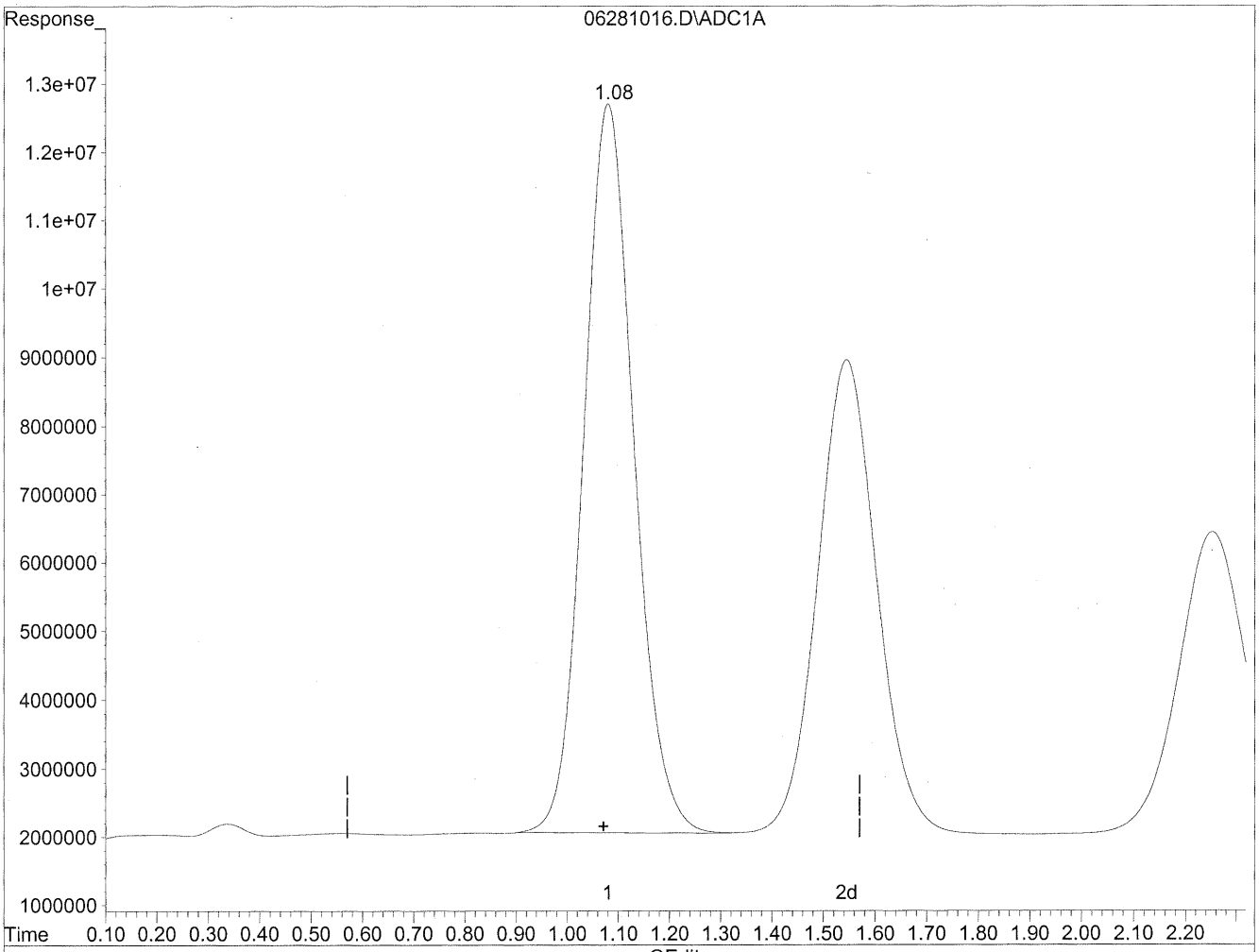
(+) = Expected Retention Time

06281002.D TO110610.M Tue Jun 29 12:58:41 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281016.D Vial: 16
Acq On : 28 Jun 2010 13:43 Operator: MD
Sample : MID CCV 5000ng/ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 28 13:57 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 5075.040ng/ml

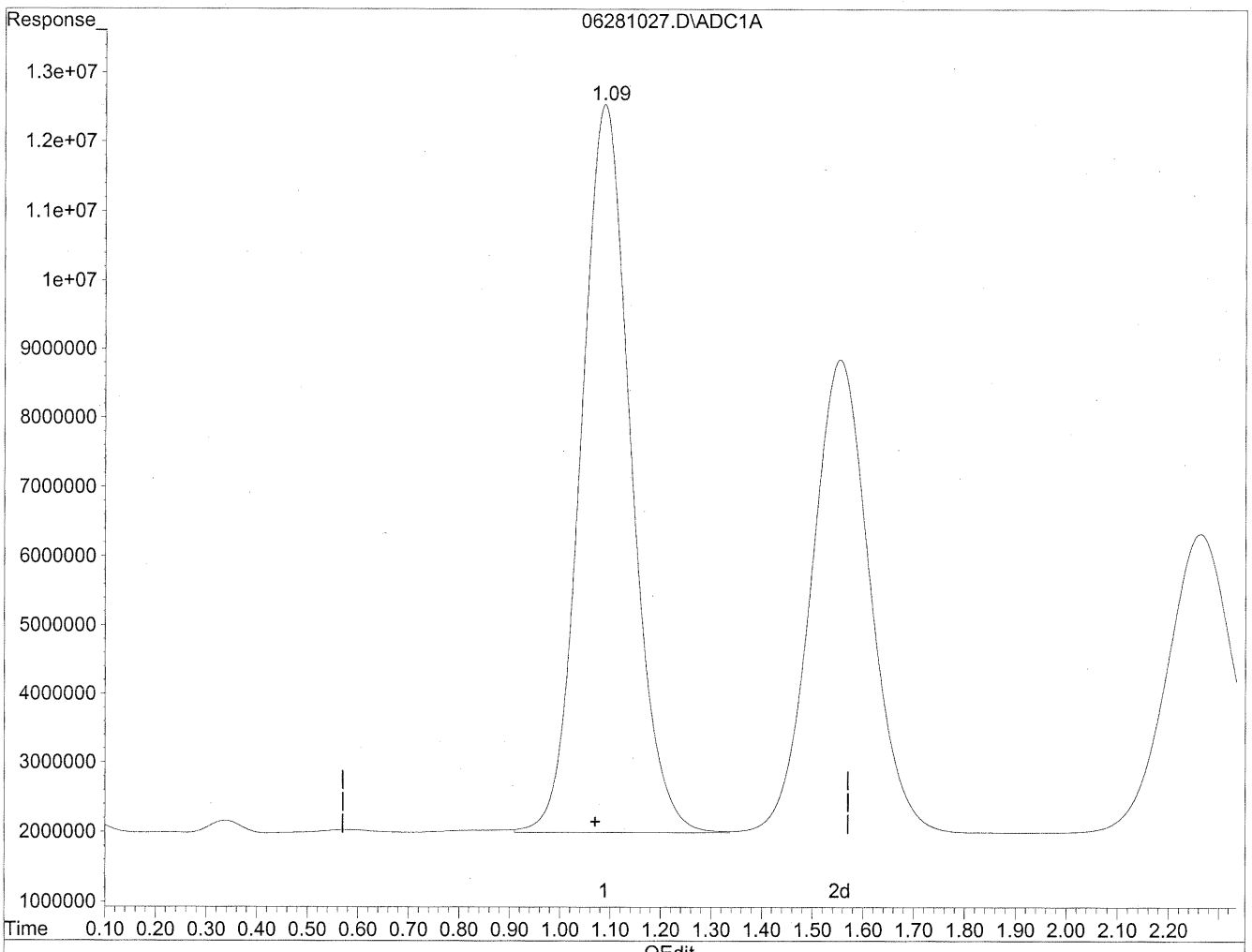
response 726672926

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281027.D Vial: 27
Acq On : 28 Jun 2010 16:12 Operator: MD
Sample : 5000ng/ml TO-11A S21-03091001 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 28 16:29 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.09min 5069.138ng/ml

response 725827950

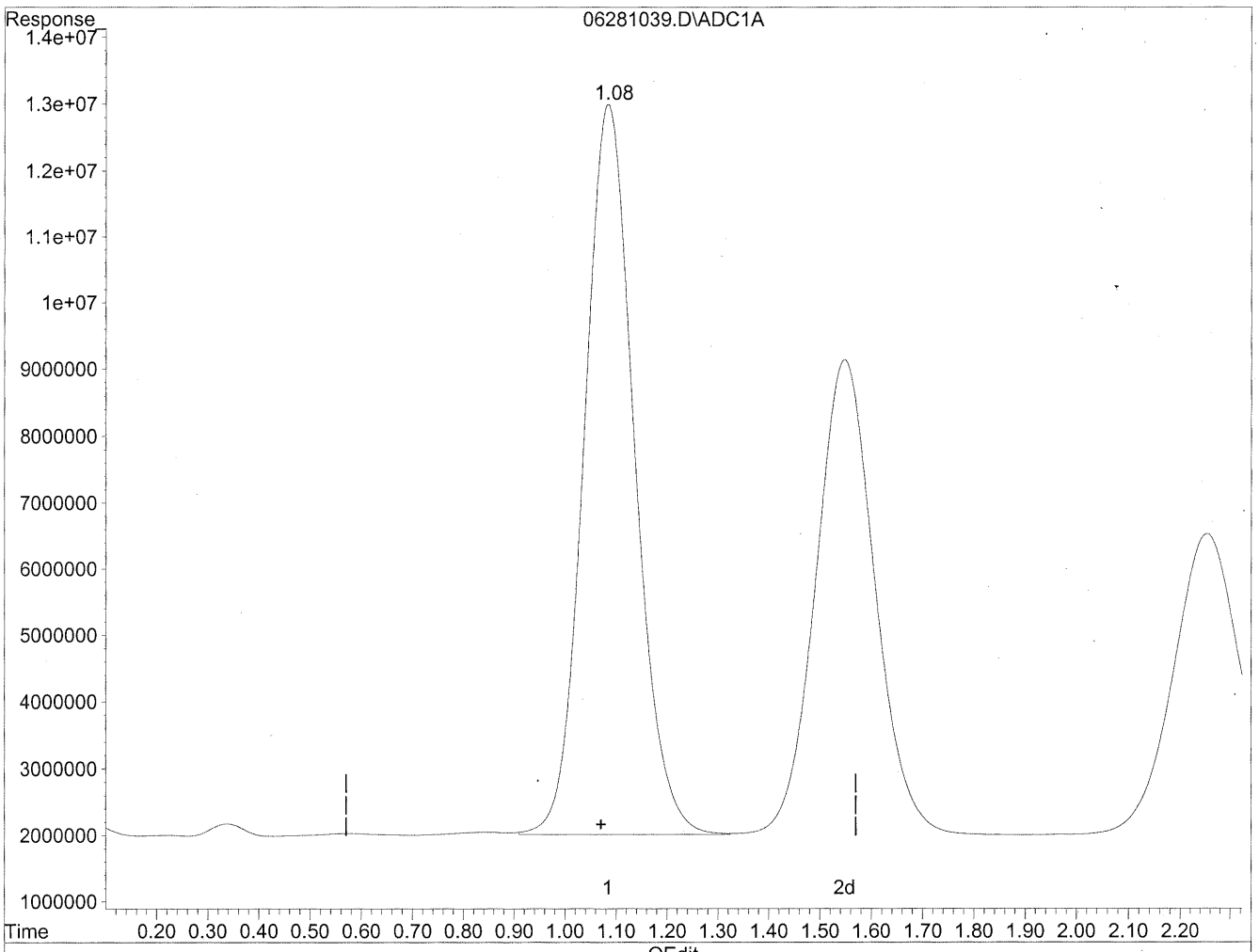
215

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281039.D Vial: 39
Acq On : 28 Jun 2010 18:55 Operator: MD
Sample : MID CCV 5000ng/ml Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 7:50 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.08min 5267.289ng/ml

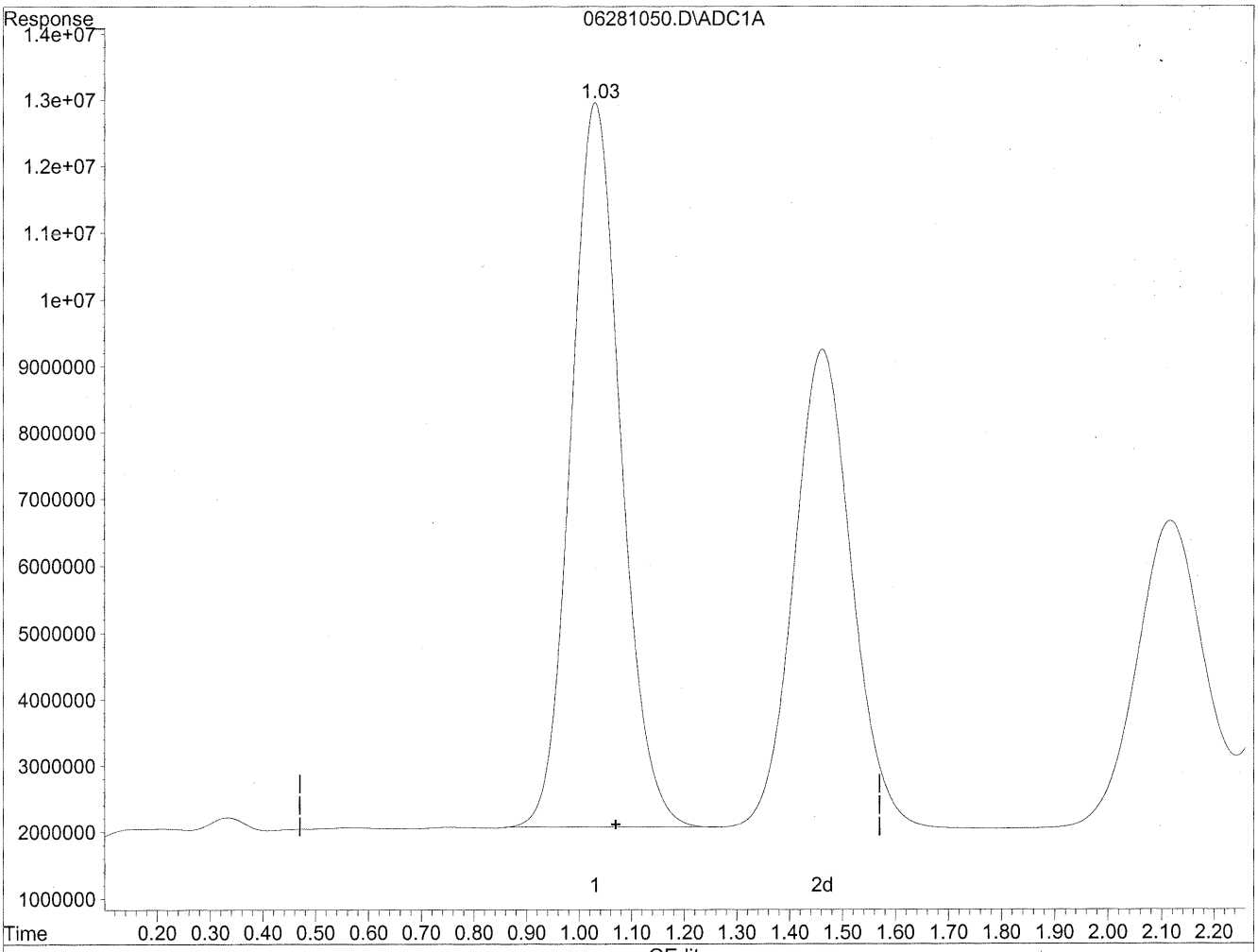
response 754200271

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281050.D Vial: 50
Acq On : 28 Jun 2010 21:24 Operator: MD
Sample : 5000ng/ml TO-11A CCV S21-03091001 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 14:58 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.03min 5021.161ng/ml

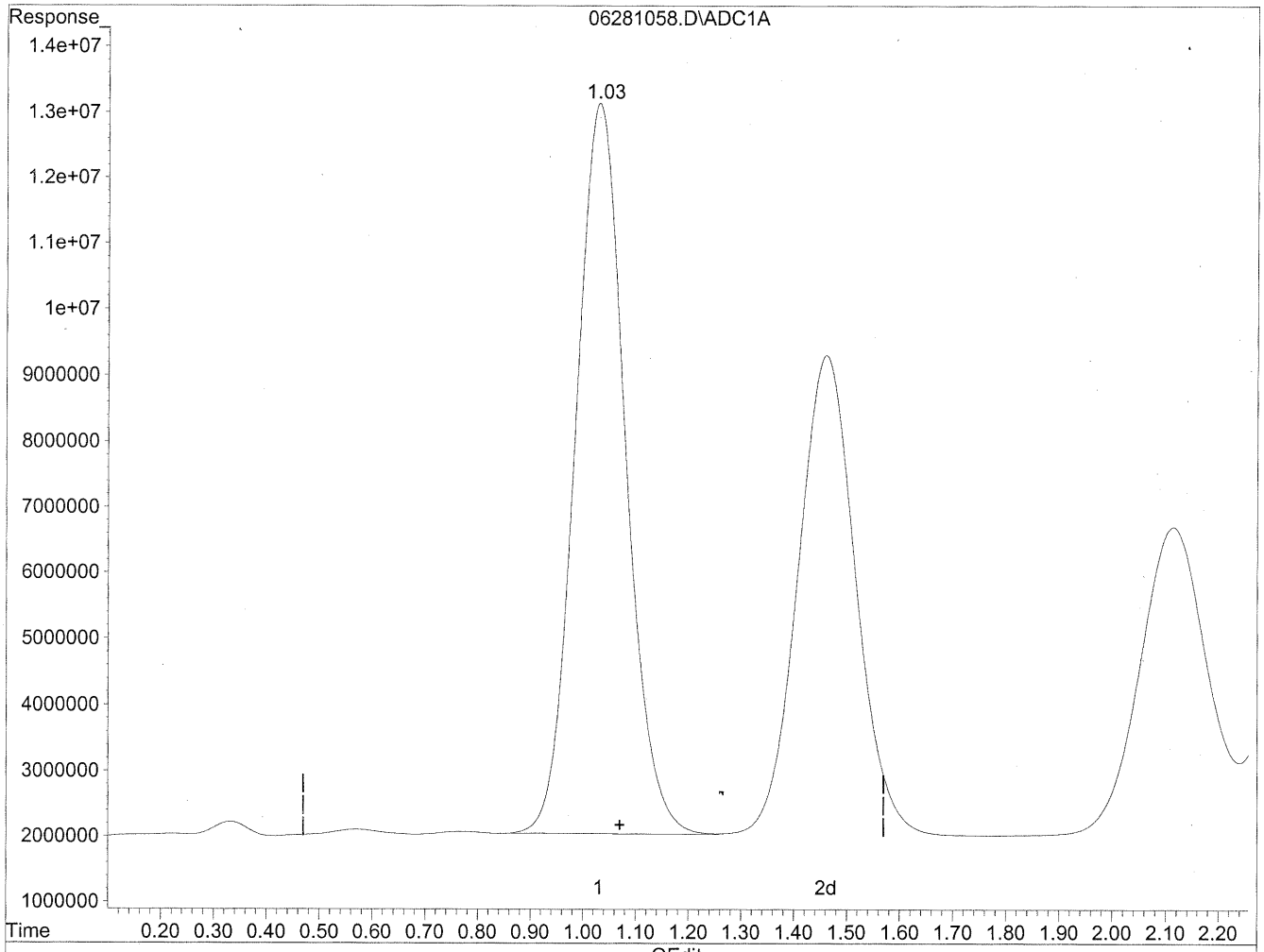
response 718958201

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\28\06281058.D Vial: 58
Acq On : 28 Jun 2010 23:12 Operator: MD
Sample : 5000ng/ml end std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:07 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.03min 5129.652ng/ml

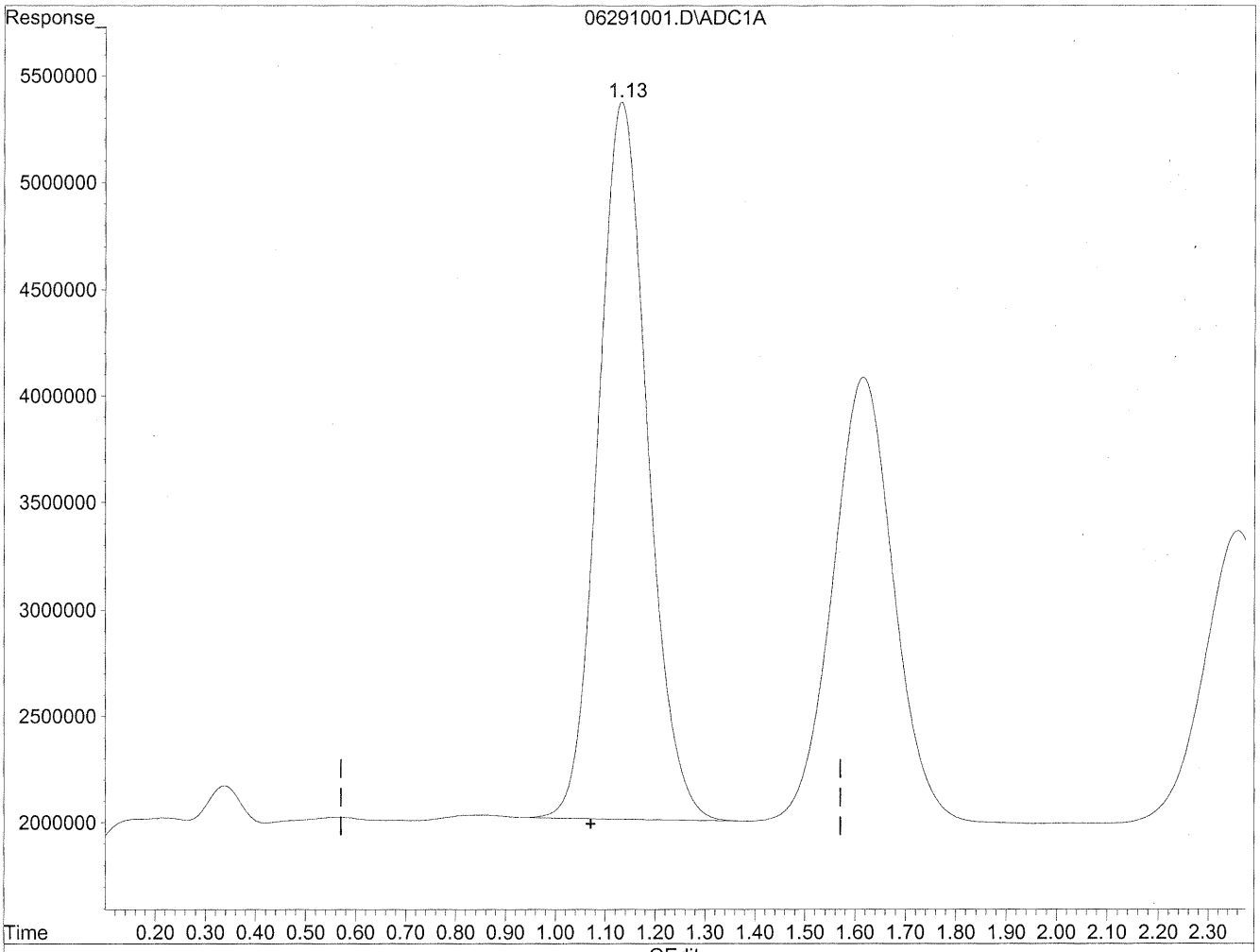
response 734492573

(+) = Expected Retention Time

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291001.D Vial: 1
Acq On : 29 Jun 2010 9:52 Operator: MD
Sample : 1500ng/ml TO-11A S21-04211003 Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 10:08 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.13min 1620.809ng/ml

response 232076587

HC 6/30/10

6/30/10

219

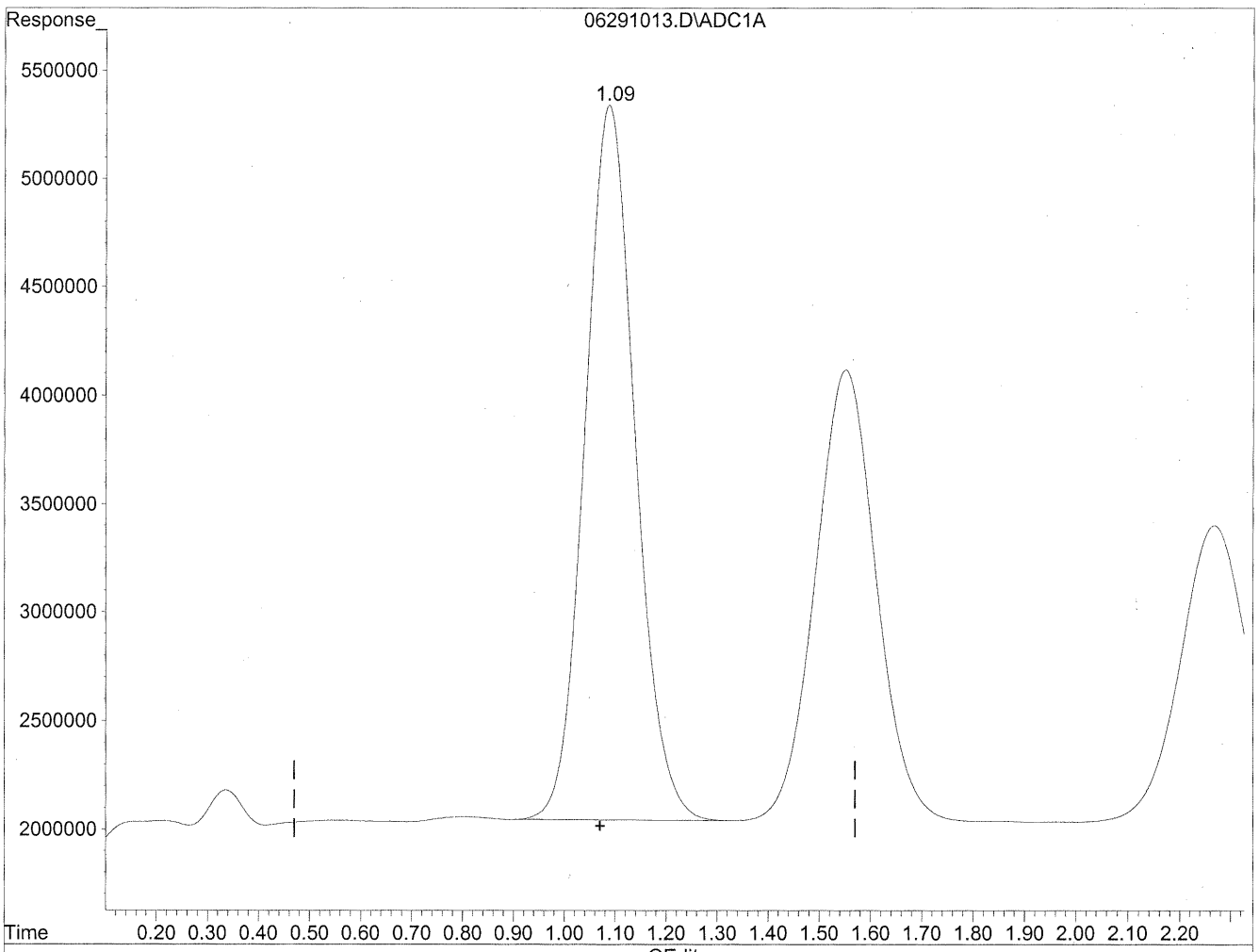
(+) = Expected Retention Time

06291001.D TO110610.M Tue Jun 29 15:07:56 2010

Quantitation Report

Data File : J:\LC01\DATA\TO11\2010_06\29\06291013.D Vial: 13
Acq On : 29 Jun 2010 12:34 Operator: MD
Sample : 1500ng/ml end std Inst : LC 01
Misc : Multiplr: 1.00
IntFile : autoint1.e
Quant Time: Jun 29 15:10 19110 Quant Results File: TO110610.RES

Method : J:\LC01\METHODS\TO110610.M (Chemstation Integrator)
Title : TO-11A Method for Aldehydes/Ketones by HPLC
Last Update : Tue Jun 29 12:58:25 2010
Response via : Multiple Level Calibration



(1) Formaldehyde

1.09min 1580.912ng/ml

response 226363940

(+) = Expected Retention Time

RUN LOGS

Injection Log

Directory: j:\lc01\data\to11\2010_06\24

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	1	06241001.d	1.	TO-11A TEST		24 Jun 110 13::
2	1	06241002.d	1.	TO-11A TEST 1.5ML/MIN		24 Jun 110 13::
3	1	06241003.d	1.	ACN blank		24 Jun 110 13::
4	2	06241004.d	1.	50ng/ml TO-11A S21-03091012		24 Jun 110 13::
5	3	06241005.d	1.	50ng/ml TO-11A S21-03091012		24 Jun 110 13::
6	4	06241006.d	1.	50ng/ml TO-11A S21-03091012		24 Jun 110 13::
7	5	06241007.d	1.	100ng/ml TO-11A S21-03091009		24 Jun 110 13::
8	6	06241008.d	1.	100ng/ml TO-11A S21-03091009		24 Jun 110 13::
9	7	06241009.d	1.	100ng/ml TO-11A S21-03091009		24 Jun 110 12::
10	8	06241010.d	1.	500ng/ml TO-11A S21-03091008		24 Jun 110 12::
11	9	06241011.d	1.	500ng/ml TO-11A S21-03091008		24 Jun 110 12::
12	10	06241012.d	1.	500ng/ml TO-11A S21-03091008		24 Jun 110 12::
13	11	06241013.d	1.	1500ng/ml TO-11A S21-04211003		24 Jun 110 12::
14	12	06241014.d	1.	1500ng/ml TO-11A S21-04211003		24 Jun 110 12::
15	13	06241015.d	1.	1500ng/ml TO-11A S21-04211003		24 Jun 110 12::
16	14	06241016.d	1.	5000ng/ml TO-11A S21-03091001		24 Jun 110 12::
17	15	06241017.d	1.	5000ng/ml TO-11A S21-03091001		24 Jun 110 12::
18	16	06241018.d	1.	5000ng/ml TO-11A S21-03091001		24 Jun 110 12::
19	17	06241019.d	1.	10000ng/ml TO-11A S21-03091007		24 Jun 110 12::
20	18	06241020.d	1.	10000ng/ml TO-11A S21-03091007		24 Jun 110 12::
21	19	06241021.d	1.	10000ng/ml TO-11A S21-03091007		24 Jun 110 12::
22	20	06241022.d	1.	~1500ng/ml TO-11A ICV S21-03091003		24 Jun 110 12::

Injection Log

Directory: j:\lc01\data\to11\2010_06\28

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	1	06281001.d	1.	prime		28 Jun 110 13::
2	2	06281002.d	1.	5000ng/ml TO-11A S21-030910001		28 Jun 110 13::
3	3	06281003.d	1.	ACN blank lot 57196HK		28 Jun 110 13::
4	4	06281004.d	1.	MB 2.0ml Lot 09240	Radiello passive sampler	28 Jun 110 13::
5	5	06281005.d	1.	P1002127-005 2.0ml		28 Jun 110 13::
6	6	06281006.d	1.	P1002127-010 2.0ml		28 Jun 110 13::
7	7	06281007.d	1.	P1002127-015 2.0ml		28 Jun 110 13::
8	8	06281008.d	1.	P1002127-020 2.0ml		28 Jun 110 13::
9	9	06281009.d	1.	P1002127-025 2.0ml		28 Jun 110 13::
10	10	06281010.d	1.	P1002127-030 2.0ml		28 Jun 110 13::
11	11	06281011.d	1.	P1002127-001 2.0ml		28 Jun 110 13::
12	12	06281012.d	1.	P1002127-001 2.0ml dup		28 Jun 110 13::
13	13	06281013.d	1.	P1002127-002 2.0ml		28 Jun 110 13::
14	14	06281014.d	1.	P1002127-003 2.0ml		28 Jun 110 12::
15	15	06281015.d	1.	ACN blank		28 Jun 110 12::
16	16	06281016.d	1.	MID CCV 5000ng/ml		28 Jun 110 12::
17	17	06281017.d	1.	P1002127-004 2.0ml		28 Jun 110 12::
18	18	06281018.d	1.	P1002127-006 2.0ml		28 Jun 110 12::
19	19	06281019.d	1.	P1002127-007 2.0ml		28 Jun 110 12::
20	20	06281020.d	1.	P1002127-008 2.0ml		28 Jun 110 12::
21	21	06281021.d	1.	P1002127-009 2.0ml		28 Jun 110 12::
22	22	06281022.d	1.	P1002127-011 2.0ml		28 Jun 110 12::
23	23	06281023.d	1.	P1002127-012 2.0ml		28 Jun 110 12::
24	24	06281024.d	1.	P1002127-013 2.0ml		28 Jun 110 12::
25	25	06281025.d	1.	P1002127-014 2.0ml		28 Jun 110 12::
26	26	06281026.d	1.	P1002127-016 2.0ml		28 Jun 110 12::
27	27	06281027.d	1.	5000ng/ml TO-11A S21-03091001		28 Jun 110 12::
28	28	06281028.d	1.	ACN blank lot 57196HK		28 Jun 110 12::
29	29	06281029.d	1.	P1002127-017 2.0ml		28 Jun 110 12::
30	30	06281030.d	1.	P1002127-017 2.0ml dup		28 Jun 110 12::
31	31	06281031.d	1.	P1002127-018 2.0ml		28 Jun 110 12::
32	32	06281032.d	1.	P1002127-019 2.0ml		28 Jun 110 12::
33	33	06281033.d	1.	P1002127-021 2.0ml		28 Jun 110 12::
34	34	06281034.d	1.	P1002127-022 2.0ml		28 Jun 110 12::
35	35	06281035.d	1.	P1002127-023 2.0ml		28 Jun 110 12::
36	36	06281036.d	1.	P1002127-024 2.0ml		28 Jun 110 12::
37	37	06281037.d	1.	P1002127-026 2.0ml		28 Jun 110 12::
38	38	06281038.d	1.	P1002127-027 2.0ml		28 Jun 110 12::
39	39	06281039.d	1.	MID CCV 5000ng/ml		28 Jun 110 12::
40	40	06281040.d	1.	P1002127-028 2.0ml		28 Jun 110 12::
41	41	06281041.d	1.	P1002127-029 2.0ml		28 Jun 110 12::
42	42	06281042.d	1.	P1002127-001 2.0ml 10x dil		28 Jun 110 12::
43	43	06281043.d	1.	P1002127-002 2.0ml 10x dil		28 Jun 110 12::
44	44	06281044.d	1.	P1002127-003 2.0ml 10x dil		28 Jun 110 12::
45	45	06281045.d	1.	P1002127-006 2.0ml 10x dil		28 Jun 110 12::
46	46	06281046.d	1.	P1002127-007 2.0ml 10x dil		28 Jun 110 12::
47	47	06281047.d	1.	P1002127-008 2.0ml 10x dil		28 Jun 110 12::
48	48	06281048.d	1.	P1002127-011 2.0ml 10x dil		28 Jun 110 12::
49	49	06281049.d	1.	P1002127-012 2.0ml 10x dil		28 Jun 110 12::
50	50	06281050.d	1.	5000ng/ml TO-11A CCV S21-03091001		28 Jun 110 12::
51	51	06281051.d	1.	ACN blank lot 57196HK		28 Jun 110 12::
52	52	06281052.d	1.	P1002127-013 2.0ml 10x dil		28 Jun 110 12::
53	53	06281053.d	1.	MB Lot 09240	Radiello passive sampler	28 Jun 110 13::
54	54	06281054.d	1.	P1002161-001 2.0ml		28 Jun 110 13::
55	55	06281055.d	1.	P1002161-002 2.0ml		28 Jun 110 13::
56	56	06281056.d	1.	P1002161-003 2.0ml		28 Jun 110 13::
57	57	06281057.d	1.	P1002161-003 2.0ml dup		28 Jun 110 13::

Injection Log

Directory: j:\lc01\data\to11\2010_06\28

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
58	58	06281058.d	1.	5000ng/ml end std		28 Jun 110 13::

Injection Log

Directory: j:\lc01\data\to11\2010_06\29

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	1	06291001.d	1.	1500ng/ml TO-11A S21-04211003		29 Jun 110 12::!
2	2	06291002.d	1.	ACN blank lot 57196HK		29 Jun 110 13::(
3	3	06291003.d	1.	P1002127-016 2.0ml 10x dil	Radiello passive sampler	29 Jun 110 13::'
4	4	06291004.d	1.	P1002127-017 2.0ml 10x dil		29 Jun 110 13::3
5	5	06291005.d	1.	P1002127-018 2.0ml 10x dil		29 Jun 110 13::4
6	6	06291006.d	1.	P1002127-021 2.0ml 10x dil		29 Jun 110 13::5
7	7	06291007.d	1.	P1002127-022 2.0ml 10x dil		29 Jun 110 13::'
8	8	06291008.d	1.	P1002127-023 2.0ml 10x dil		29 Jun 110 13::2
9	9	06291009.d	1.	P1002127-026 2.0ml 10x dil		29 Jun 110 13::4
10	10	06291010.d	1.	P1002127-027 2.0ml 10x dil		29 Jun 110 13::5
11	11	06291011.d	1.	P1002127-028 2.0ml 10x dil		29 Jun 110 13::(
12	12	06291012.d	1.	P1002161-003 2.0ml 10x dil		29 Jun 110 13::2
13	13	06291013.d	1.	1500ng/ml end std		29 Jun 110 13::3