

**Air
Toxics LTD.**
Laboratory Services Since 1989

Electronic Comprehensive Validation Package (eCVP)



AN ENVIRONMENTAL ANALYTICAL LABORATORY

COMPREHENSIVE VALIDATION PACKAGE

Modified TO-11A

INVENTORY SHEET

Work Order #: 1011114B

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Comments:

Completed by:

Vera Belitsky

(Signature)

Vera Belitsky / Document Control

(Print Name & Title)

11/19/10

(Date)

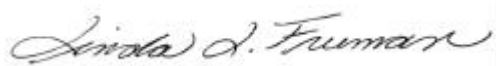
WORK ORDER #: 1011114B

Work Order Summary

CLIENT:	Mr. Brian Baker Environmental Health & Engineering, Inc. 117 Fourth Avenue Needham, MA 02494	BILL TO:	Accounts Payable Environmental Health & Engineering, Inc. 117 Fourth Avenue Needham, MA 02494
PHONE:	800-825-5343	P.O. #	17131
FAX:	781-247-4305	PROJECT #	17131
DATE RECEIVED:	11/04/2010	CONTACT:	Ausha Scott
DATE COMPLETED:	11/17/2010		

<u>FRACTION #</u>	<u>NAME</u>	<u>TEST</u>
16A	118635	Modified TO-11A
17A	118636	Modified TO-11A
18A	118637	Modified TO-11A
19A	118638	Modified TO-11A
20A	118639	Modified TO-11A
21A	118651	Modified TO-11A
22A	118652	Modified TO-11A
23A	118653	Modified TO-11A
24A	118654	Modified TO-11A
25A	118655	Modified TO-11A
26A	Lab Blank	Modified TO-11A
27A	LCS	Modified TO-11A
27AA	LCSD	Modified TO-11A

CERTIFIED BY:



Laboratory Director

DATE: 11/18/10

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LABORATORY NARRATIVE
Modified TO-11A
Environmental Health & Engineering, Inc.
Workorder# 101114B

Ten Radiello 165 (Aldehydes) samples were received on November 04, 2010. The laboratory performed analysis via modified Method TO-11A using reverse phase High Pressure Liquid Chromatography (HPLC) with an Ultraviolet (UV) Detector. The method involves eluting the sorbent tubes with acetonitrile using a gravity feed technique. Method modifications taken to run these samples include:

Requirement	TO-11A	ATL Modifications
ACN Purity Check	Contribution of analytes from ACN determined as described Sections 9.1.1 and 9.1.2 of Compendium TO-11A.	Total contribution of analytes from ACN and cartridge combined is determined.
Initial Calibration Curve (ICAL)	Multi-point using linear regression performed every 6 months; $r^2 > 0.999$	Multi-point using average Response Factor; % RSD ≤ 10 %. Re-calibration if daily cal. fails, major maintenance, or column change. Linear regression is performed when requested.
Blank Subtraction	Average blank concentrations calculated. Blank value subtracted from sample result.	One Lab Blank is analyzed per batch; no blank subtraction performed on samples.

Receiving Notes

There were no receiving discrepancies.

Analytical Notes

The recovery for Formaldehyde in the LCS was outside the laboratory control limits. Due to the nature of the media, it is not possible to re-extract the associated samples.

Definition of Data Qualifying Flags

Seven qualifiers may have been used on the data analysis sheets and indicate as follows:

- B - Compound present in laboratory blank greater than reporting limit.
- J - Estimated value.
- E - Exceeds instrument calibration range.
- S - Saturated peak.
- Q - Exceeds quality control limits.
- U - Compound analyzed for but not detected above the detection limit.
- M - Reported value may be biased due to apparent matrix interferences.

File extensions may have been used on the data analysis sheets and indicates as follows:

- a-File was requantified

b-File was quantified by a second column and detector
r1-File was requantified for the purpose of reissue

Table 1

Client Sample ID	Lab Sample ID	Date Collected	Date Received	Date Extracted	Sample	Sample Extract		
					Holding Time (Days)	Date Analyzed	Holding Time (Days)	Sample Condition
118635	1011114B-16A	11/ 1/2010	11/ 4/2010	11/ 9/2010	8	11/ 9/2010	0	Good
118636	1011114B-17A	11/ 1/2010	11/ 4/2010	11/ 9/2010	8	11/ 9/2010	0	Good
118637	1011114B-18A	11/ 1/2010	11/ 4/2010	11/ 9/2010	8	11/ 9/2010	0	Good
118638	1011114B-19A	11/ 1/2010	11/ 4/2010	11/ 9/2010	8	11/ 9/2010	0	Good
118639	1011114B-20A	NA	11/ 4/2010	11/ 9/2010	NA	11/ 9/2010	0	Good
118651	1011114B-21A	11/ 1/2010	11/ 4/2010	11/ 9/2010	8	11/ 9/2010	0	Good
118652	1011114B-22A	11/ 1/2010	11/ 4/2010	11/ 9/2010	8	11/ 9/2010	0	Good
118653	1011114B-23A	11/ 1/2010	11/ 4/2010	11/ 9/2010	8	11/ 9/2010	0	Good
118654	1011114B-24A	11/ 1/2010	11/ 4/2010	11/ 9/2010	8	11/ 9/2010	0	Good
118655	1011114B-25A	NA	11/ 4/2010	11/ 9/2010	NA	11/ 9/2010	0	Good
Lab Blank	1011114B-26A	NA	NA	11/ 9/2010	NA	11/ 9/2010	0	Good
LCS	1011114B-27A	NA	NA	11/ 9/2010	NA	11/ 9/2010	0	Good
LCSD	1011114B-27AA	NA	NA	11/ 9/2010	NA	11/ 9/2010	0	Good

Sample Results and Raw Data

Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118635

Lab ID#: 1011114B-16A

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	10	45

Client Sample ID: 118635

Lab ID#: 1011114B-16A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1109038
Dil. Factor: 10.0

Date of Collection: 11/1/10
Date of Analysis: 11/9/10 09:30 PM
Date of Extraction: 11/9/10

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	10	45

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109038.d
Lab Smp Id: 1011114B-16A Client Smp ID: 10
Inj Date : 09-NOV-2010 21:30
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1011114B-16A;10
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1
Dil Factor: 10.00000
Integrator: Falcon Compound Sublist: form.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	10.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	2.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					CONCENTRATIONS	
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.861	2.957	-0.096	2115312		
2 FORMALDEHYDE	3.634	3.731	-0.097	1222341	11.1656	44.7

Data File: /chem/hp1cf.i/09Nov2010.b/f1109038.d

Date : 09-NOV-2010 21:30

Client ID: 10

Sample Info: j101114B-16A;10

Purge Volume: 5.0

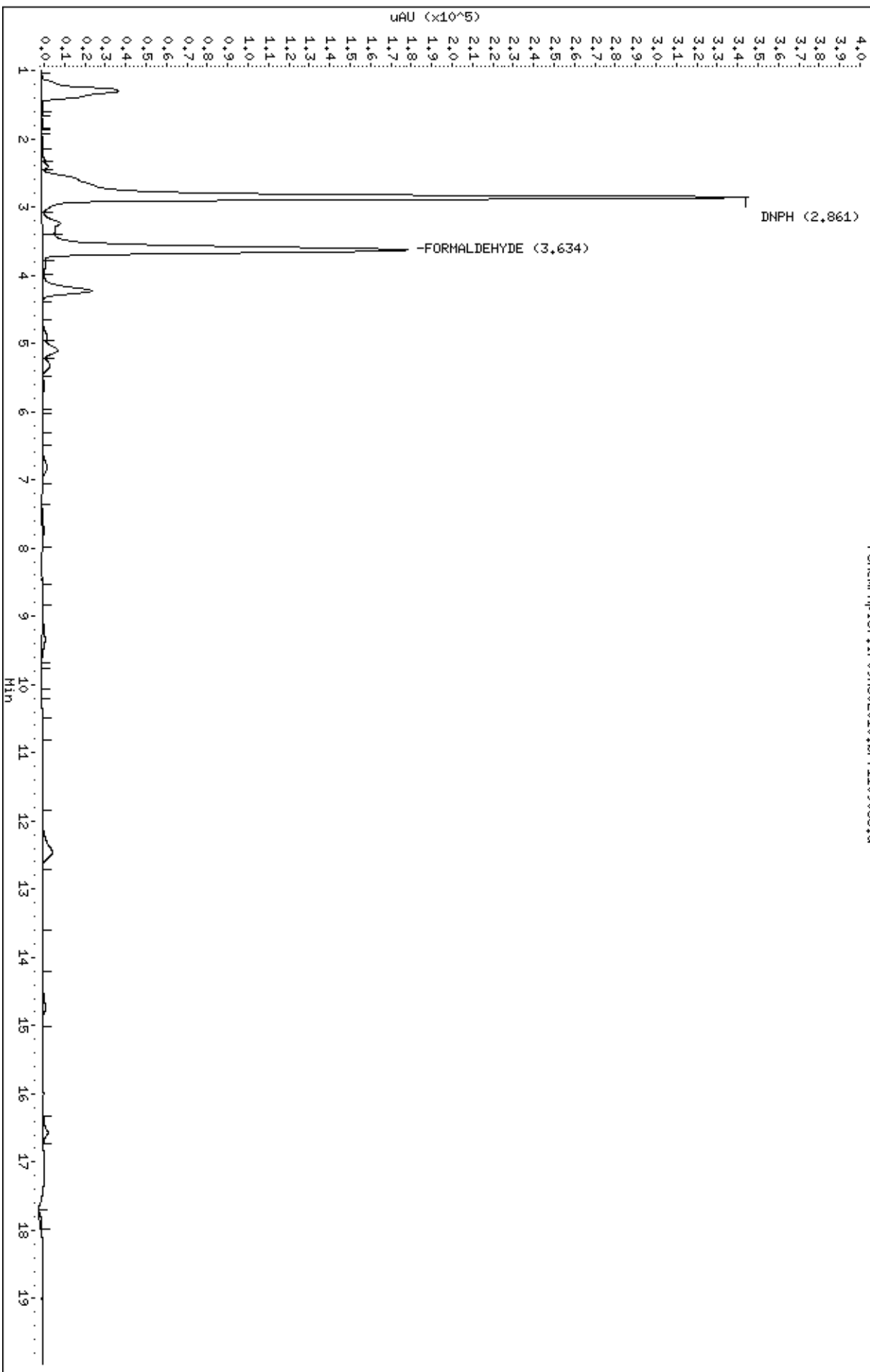
Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00

/chem/hp1cf.i/09Nov2010.b/f1109038.d



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118636

Lab ID#: 1011114B-17A

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	10	43

Client Sample ID: 118636

Lab ID#: 1011114B-17A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1109041
Dil. Factor: 10.0

Date of Collection: 11/1/10
Date of Analysis: 11/9/10 10:33 PM
Date of Extraction: 11/9/10

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	10	43

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
 Data file : /chem/hplcf.i/09Nov2010.b/f1109041.d
 Lab Smp Id: 1011114B-17A Client Smp ID: 10
 Inj Date : 09-NOV-2010 22:33
 Operator : CRL Inst ID: hplcf.i
 Smp Info : ;1011114B-17A;10
 Misc Info :
 Comment :
 Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
 Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
 Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
 Als bottle: 1
 Dil Factor: 10.00000
 Integrator: Falcon Compound Sublist: form.sub
 Target Version: 3.50
 Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	10.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	2.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	CONCENTRATIONS	
							ON-COLUMN (ng)	FINAL (ug)
1 DNPH	2.926	2.957	-0.031			1288752		
2 FORMALDEHYDE	3.733	3.731	0.002			1178047	10.7610	43.0

Data File: /chem/hp1cf.i/09Nov2010.b/f1109041.d

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Date : 09-NOV-2010 22:33

Client ID: 10

Instrument: hp1cf.i

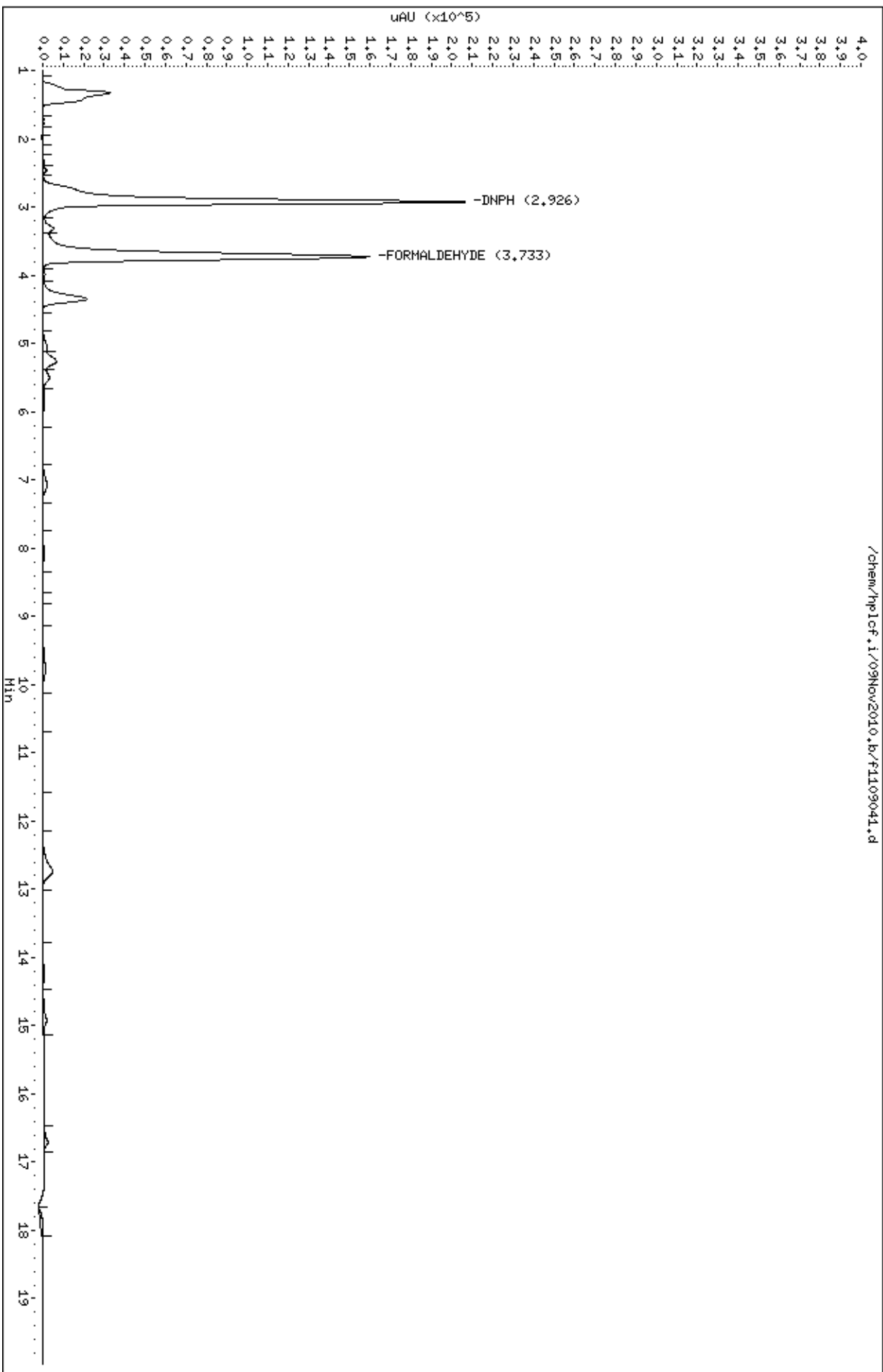
Sample Info: j1011114B-17A;10

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118637

Lab ID#: 1011114B-18A

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	10	50

Client Sample ID: 118637

Lab ID#: 1011114B-18A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1109042
Dil. Factor: 10.0

Date of Collection: 11/1/10
Date of Analysis: 11/9/10 10:54 PM
Date of Extraction: 11/9/10

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	10	50

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109042.d
Lab Smp Id: 1011114B-18A Client Smp ID: 10
Inj Date : 09-NOV-2010 22:54
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1011114B-18A;10
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1
Dil Factor: 10.00000
Integrator: Falcon Compound Sublist: form.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	10.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	2.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	CONCENTRATIONS	
							ON-COLUMN	FINAL
							(ng)	(ug)
1 DNPH	2.926	2.957	-0.031			2555609		
2 FORMALDEHYDE	3.726	3.731	-0.005			1373465	12.5460	50.2

Data File: /chem/hplcf.i/09Nov2010.b/f1109042.d

Date : 09-NOV-2010 22:54

Client ID: 10

Sample Info: j1011114B-18A;10

Purge Volume: 5.0

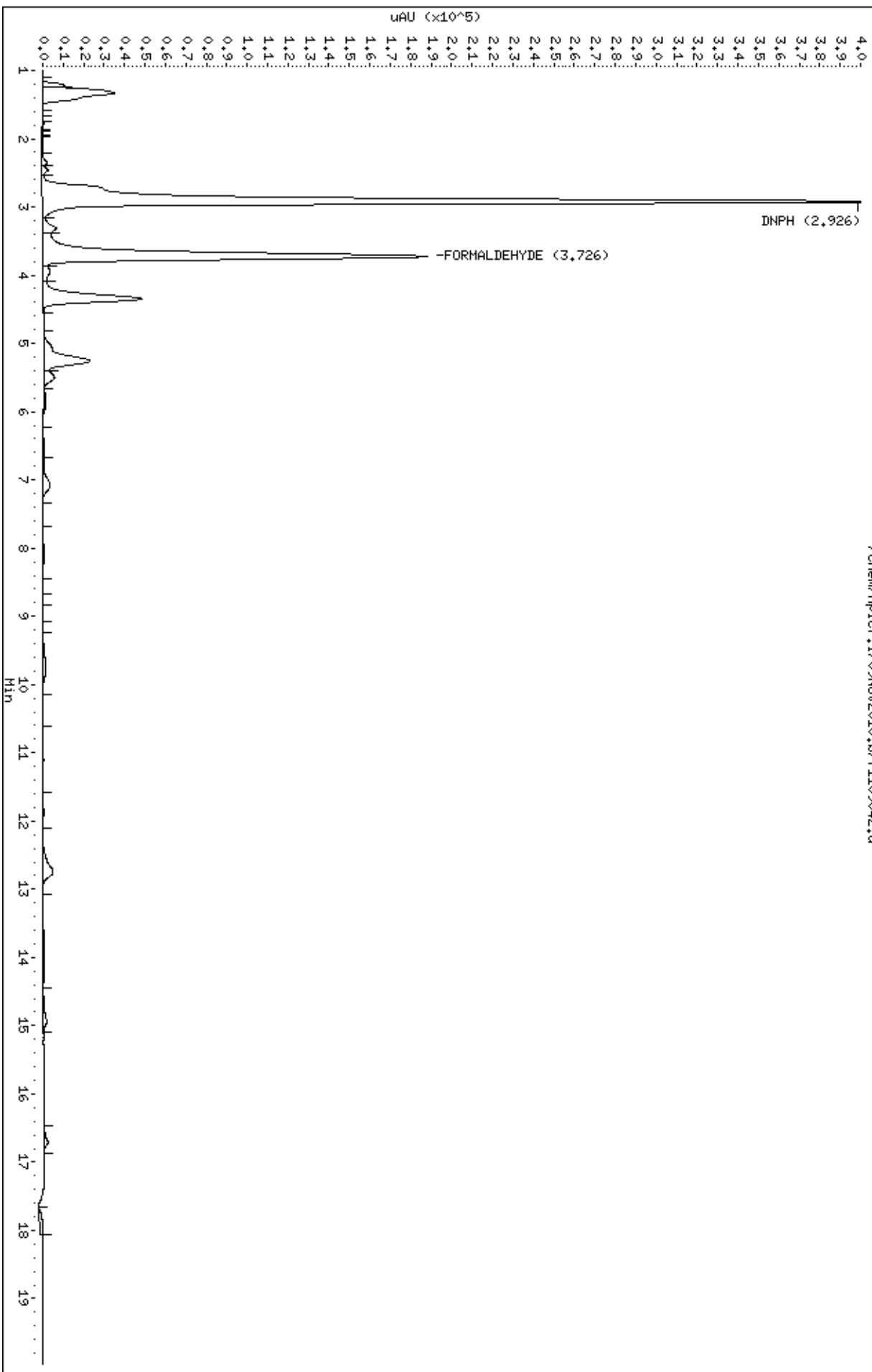
Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: CRL

Column diameter: 2.00

/chem/hplcf.i/09Nov2010.b/f1109042.d



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118638

Lab ID#: 1011114B-19A

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	1.0	2.1

Client Sample ID: 118638

Lab ID#: 1011114B-19A

RADIELLO ANALYSIS BY HPLC/UV

File Name:	f1109025	Date of Collection: 11/1/10
Dil. Factor:	1.00	Date of Analysis: 11/9/10 04:59 PM
		Date of Extraction: 11/9/10

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	1.0	2.1

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109025.d
Lab Smp Id: 1011114B-19A
Inj Date : 09-NOV-2010 16:59
Operator : CRL
Smp Info : ;1011114B-19A;
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf
Cal Date : 14-SEP-2010 18:30
Als bottle: 1
Dil Factor: 1.00000
Integrator: Falcon
Target Version: 3.50
Processing Host: eeyore

Inst ID: hplcf.i

Quant Type: ESTD

Cal File: f0914014.d

Compound Sublist: form.sub

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	2.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds					CONCENTRATIONS	
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.932	2.957	-0.025	19199775		
2 FORMALDEHYDE	3.732	3.731	0.001	574094	5.24410	2.10

Data File: /chem/hp1cf.i/09Nov2010.b/f1109025.d

Date : 09-NOV-2010 16:59

Client ID:

Sample Info: j101114B-19A;

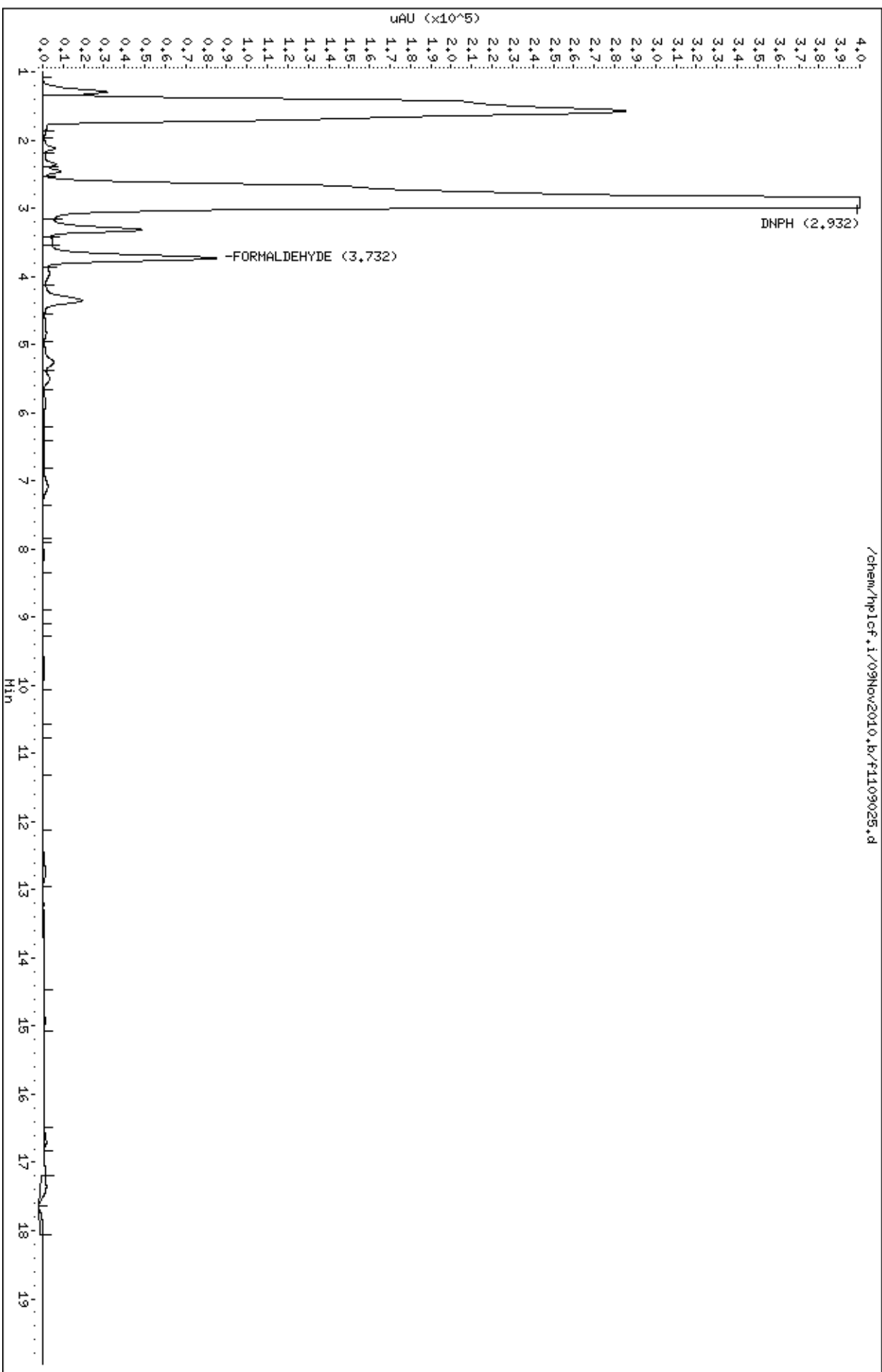
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



**Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV**

Client Sample ID: 118639

Lab ID#: 1011114B-20A

No Detections Were Found.

Client Sample ID: 118639

Lab ID#: 1011114B-20A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1109020
Dil. Factor: 1.00

Date of Collection: NA
Date of Analysis: 11/9/10 03:14 PM
Date of Extraction: 11/9/10

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	1.0	Not Detected

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109020.d
Lab Smp Id: 1011114B-20A
Inj Date : 09-NOV-2010 15:14
Operator : CRL
Smp Info : ;1011114B-20A;
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf
Cal Date : 14-SEP-2010 18:30
Als bottle: 1
Dil Factor: 1.00000
Integrator: Falcon
Target Version: 3.50
Processing Host: eeyore

Inst ID: hplcf.i

Quant Type: ESTD

Cal File: f0914014.d

Compound Sublist: form.sub

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	2.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	CONCENTRATIONS	
							ON-COLUMN (ng)	FINAL (ug)
1 DNPH	2.925	2.957	-0.032			28645449		
2 FORMALDEHYDE						Compound Not Detected.		

Data File: /chem/hp1cf.i/09Nov2010.b/f1109020.d

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Date : 09-NOV-2010 15:14

Client ID:

Instrument: hp1cf.i

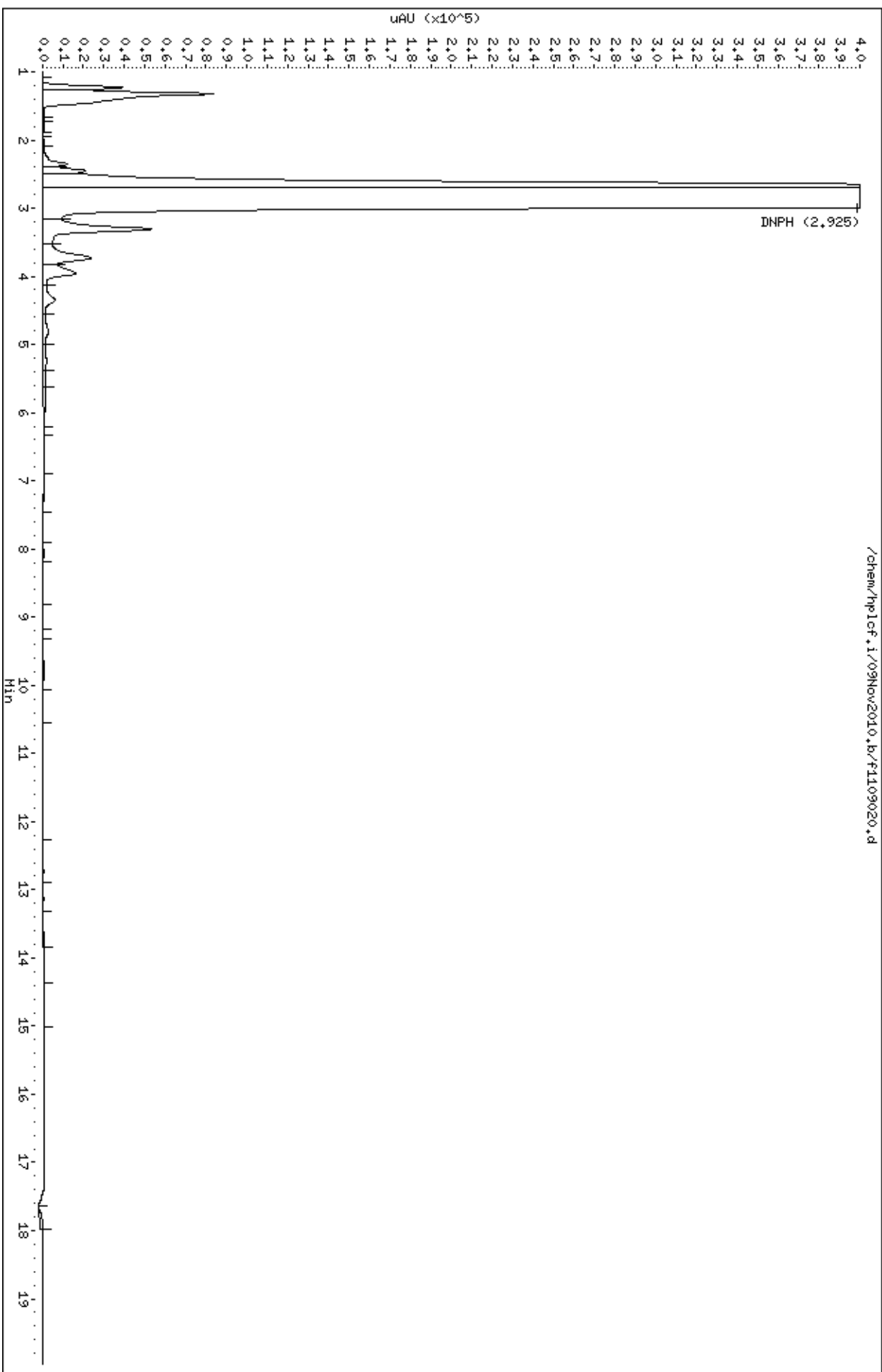
Sample Info: j101114B-20A;

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118651

Lab ID#: 1011114B-21A

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	10	81

Client Sample ID: 118651

Lab ID#: 1011114B-21A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1109043
Dil. Factor: 10.0

Date of Collection: 11/1/10
Date of Analysis: 11/9/10 11:15 PM
Date of Extraction: 11/9/10

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	10	81

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
 Data file : /chem/hplcf.i/09Nov2010.b/f1109043.d
 Lab Smp Id: 1011114B-21A Client Smp ID: 10
 Inj Date : 09-NOV-2010 23:15
 Operator : CRL Inst ID: hplcf.i
 Smp Info : ;1011114B-21A;10
 Misc Info :
 Comment :
 Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
 Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
 Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
 Als bottle: 1
 Dil Factor: 10.00000
 Integrator: Falcon Compound Sublist: form.sub
 Target Version: 3.50
 Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	10.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	2.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	CONCENTRATIONS	
							ON-COLUMN (ng)	FINAL (ug)
1 DNPH	2.930	2.957	-0.027		2145771			
2 FORMALDEHYDE	3.730	3.731	-0.001		2217716	20.2579	81.0	

Data File: /chem/hp1cf.i/09Nov2010.b/f1109043.d

Page 1

Date : 09-NOV-2010 23:15

Client ID: 10

Instrument: hp1cf.i

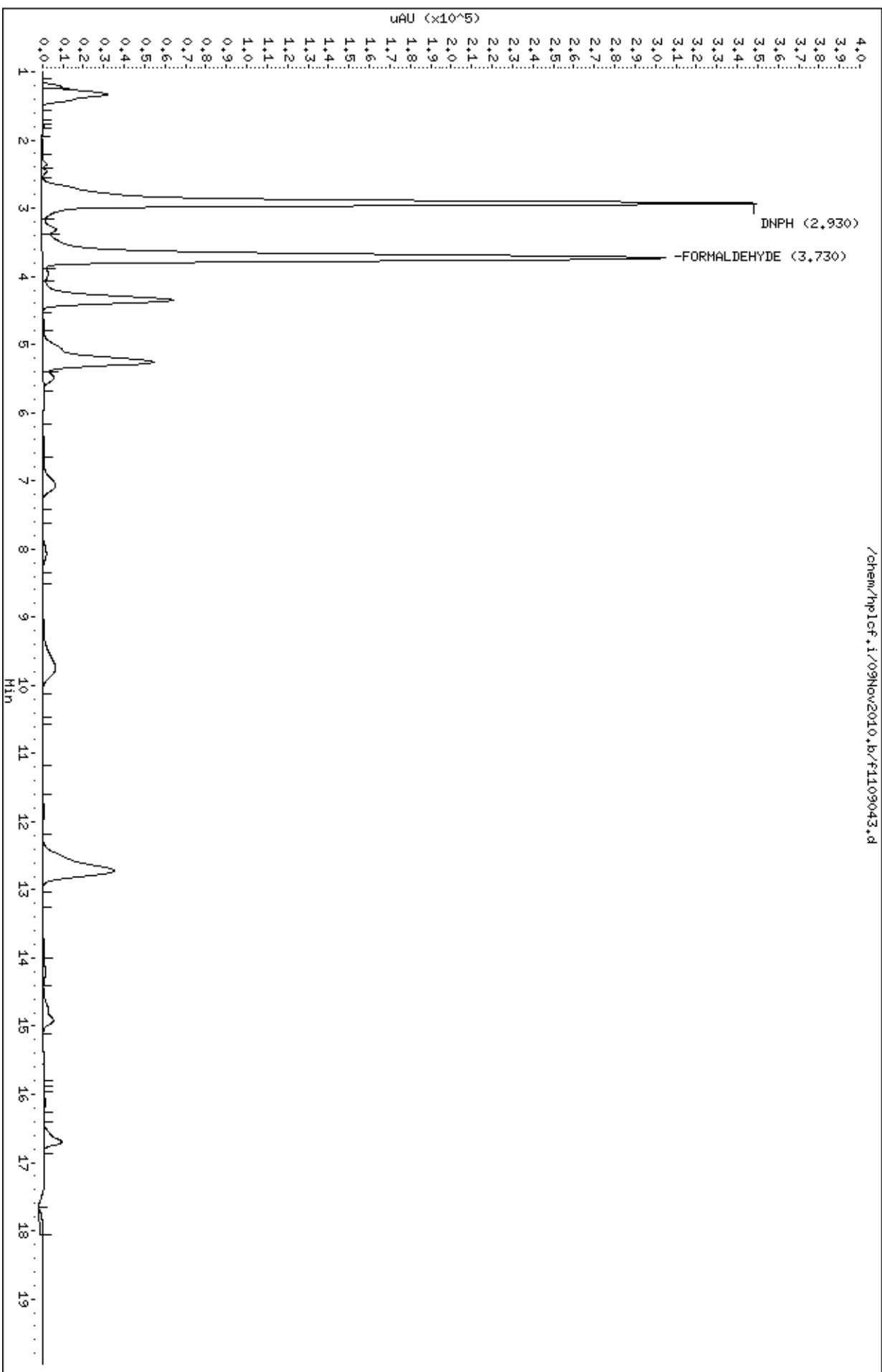
Sample Info: j1011114B-21A;10

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118652

Lab ID#: 1011114B-22A

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	10	62

Client Sample ID: 118652

Lab ID#: 1011114B-22A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1109044
Dil. Factor: 10.0

Date of Collection: 11/1/10
Date of Analysis: 11/9/10 11:35 PM
Date of Extraction: 11/9/10

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	10	62

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109044.d
Lab Smp Id: 1011114B-22A Client Smp ID: 10
Inj Date : 09-NOV-2010 23:35
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1011114B-22A;10
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1
Dil Factor: 10.00000
Integrator: Falcon Compound Sublist: form.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	10.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	2.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					CONCENTRATIONS	
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.924	2.957	-0.033	2065801		
2 FORMALDEHYDE	3.724	3.731	-0.007	1683295	15.3762	61.5

Data File: /chem/hp1cf.i/09Nov2010.b/f1109044.d

Date : 09-NOV-2010 23:35

Client ID: 10

Sample Info: j1011114B-22A;10

Purge Volume: 5.0

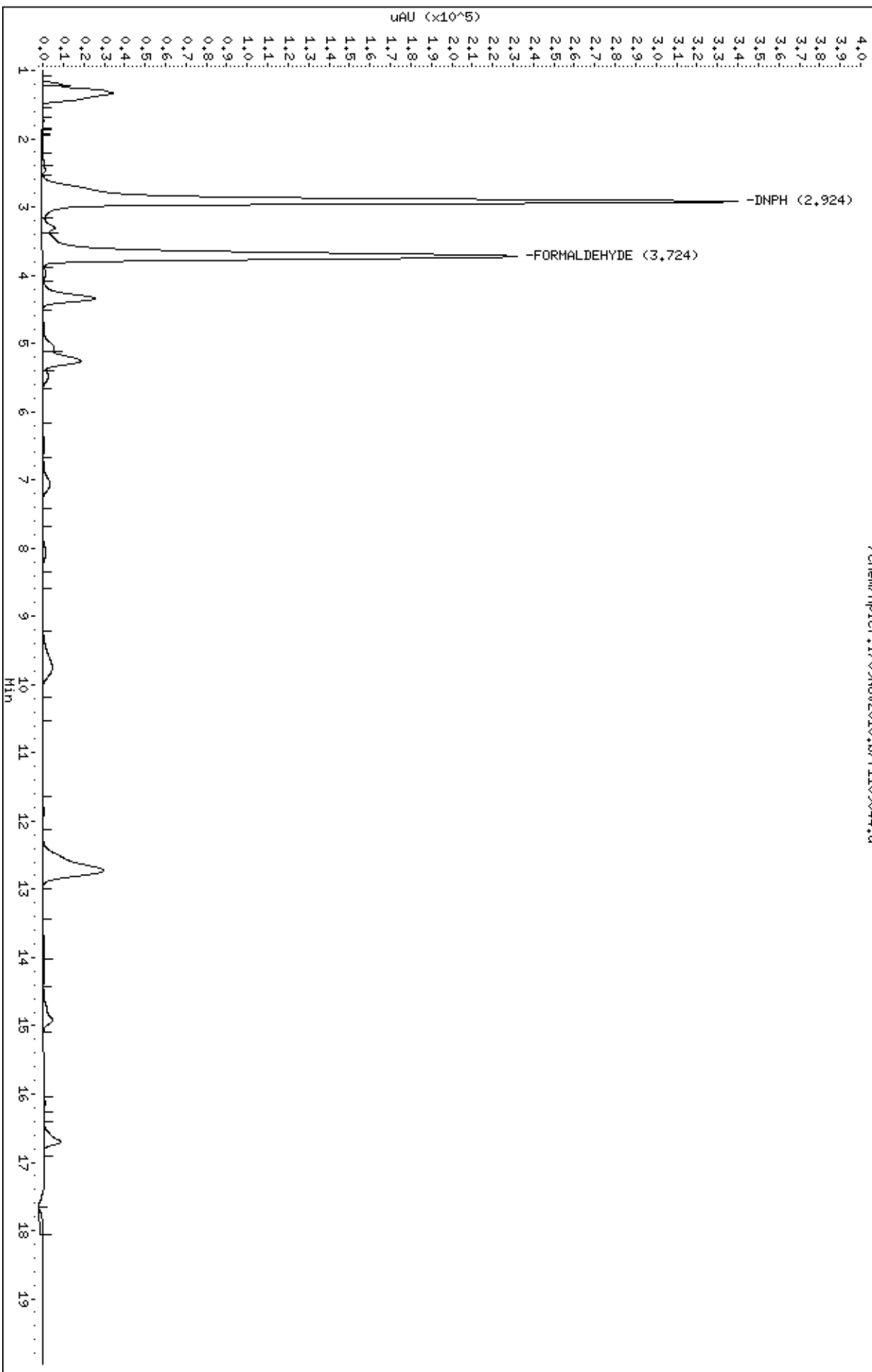
Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00

/chem/hp1cf.i/09Nov2010.b/f1109044.d



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118653

Lab ID#: 1011114B-23A

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	10	54

Client Sample ID: 118653

Lab ID#: 1011114B-23A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1109045
Dil. Factor: 10.0

Date of Collection: 11/1/10
Date of Analysis: 11/9/10 11:56 PM
Date of Extraction: 11/9/10

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	10	54

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109045.d
Lab Smp Id: 1011114B-23A Client Smp ID: 10
Inj Date : 09-NOV-2010 23:56
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1011114B-23A;10
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1
Dil Factor: 10.00000
Integrator: Falcon Compound Sublist: form.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	10.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	2.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					CONCENTRATIONS	
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.923	2.957	-0.034	1193512		
2 FORMALDEHYDE	3.723	3.731	-0.008	1466278	13.3938	53.6

Data File: /chem/hp1cf.i/09Nov2010.b/f1109045.d

Date : 09-NOV-2010 23:56

Client ID: 10

Sample Info: j1011114B-23A;10

Purge Volume: 5.0

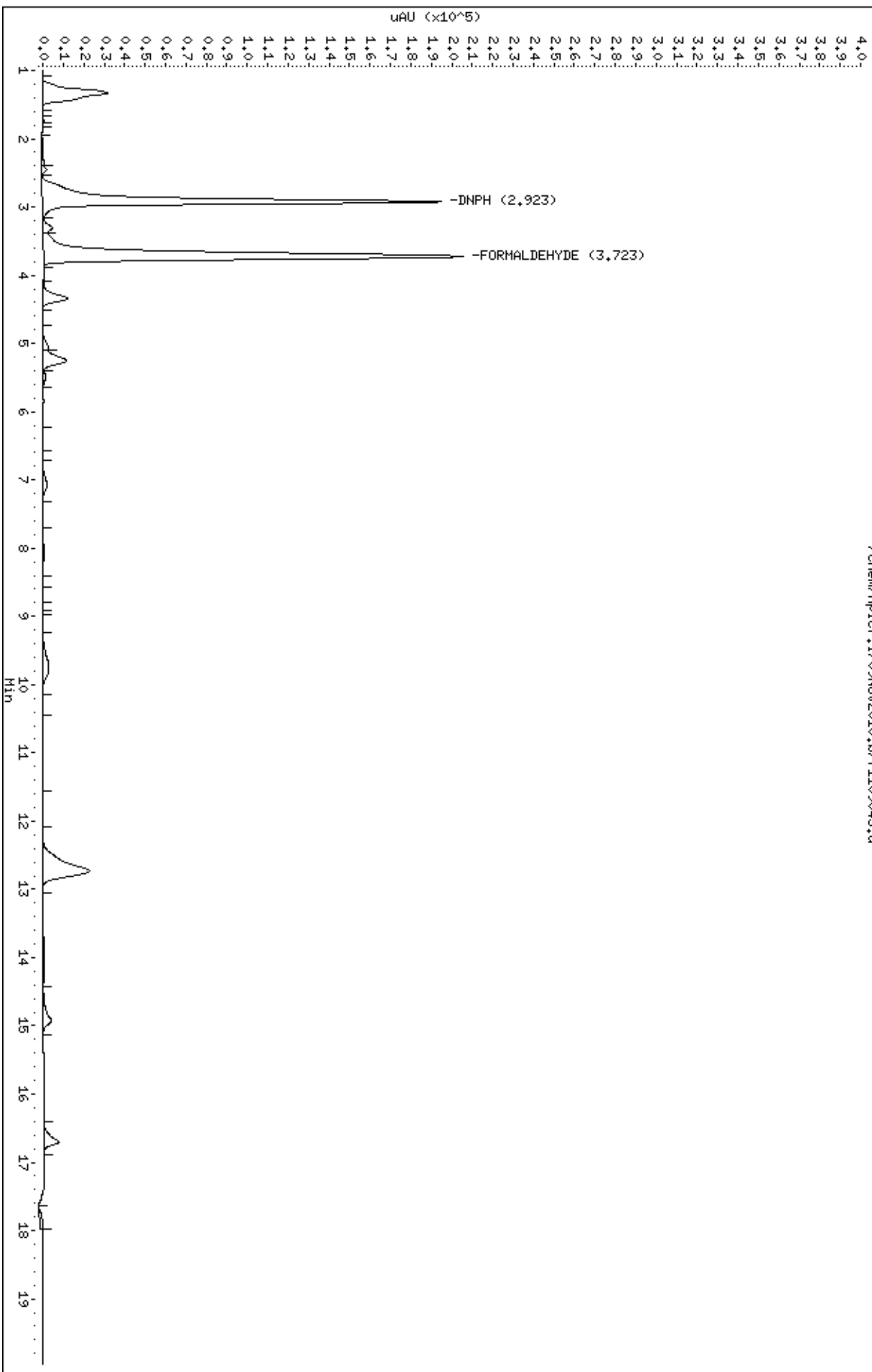
Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00

/chem/hp1cf.i/09Nov2010.b/f1109045.d



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118654

Lab ID#: 1011114B-24A

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	1.0	2.5

Client Sample ID: 118654

Lab ID#: 1011114B-24A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1109026
Dil. Factor: 1.00

Date of Collection: 11/1/10
Date of Analysis: 11/9/10 05:19 PM
Date of Extraction: 11/9/10

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	1.0	2.5

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109026.d
Lab Smp Id: 1011114B-24A
Inj Date : 09-NOV-2010 17:19
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1011114B-24A;
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: form.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	2.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					CONCENTRATIONS	
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.955	2.957	-0.002	20291783		
2 FORMALDEHYDE	3.755	3.731	0.024	692761	6.32808	2.53

Data File: /chem/hp1cf.i/09Nov2010.b/f1109026.d

Page 1

Date : 09-NOV-2010 17:19

Client ID:

Instrument: hp1cf.i

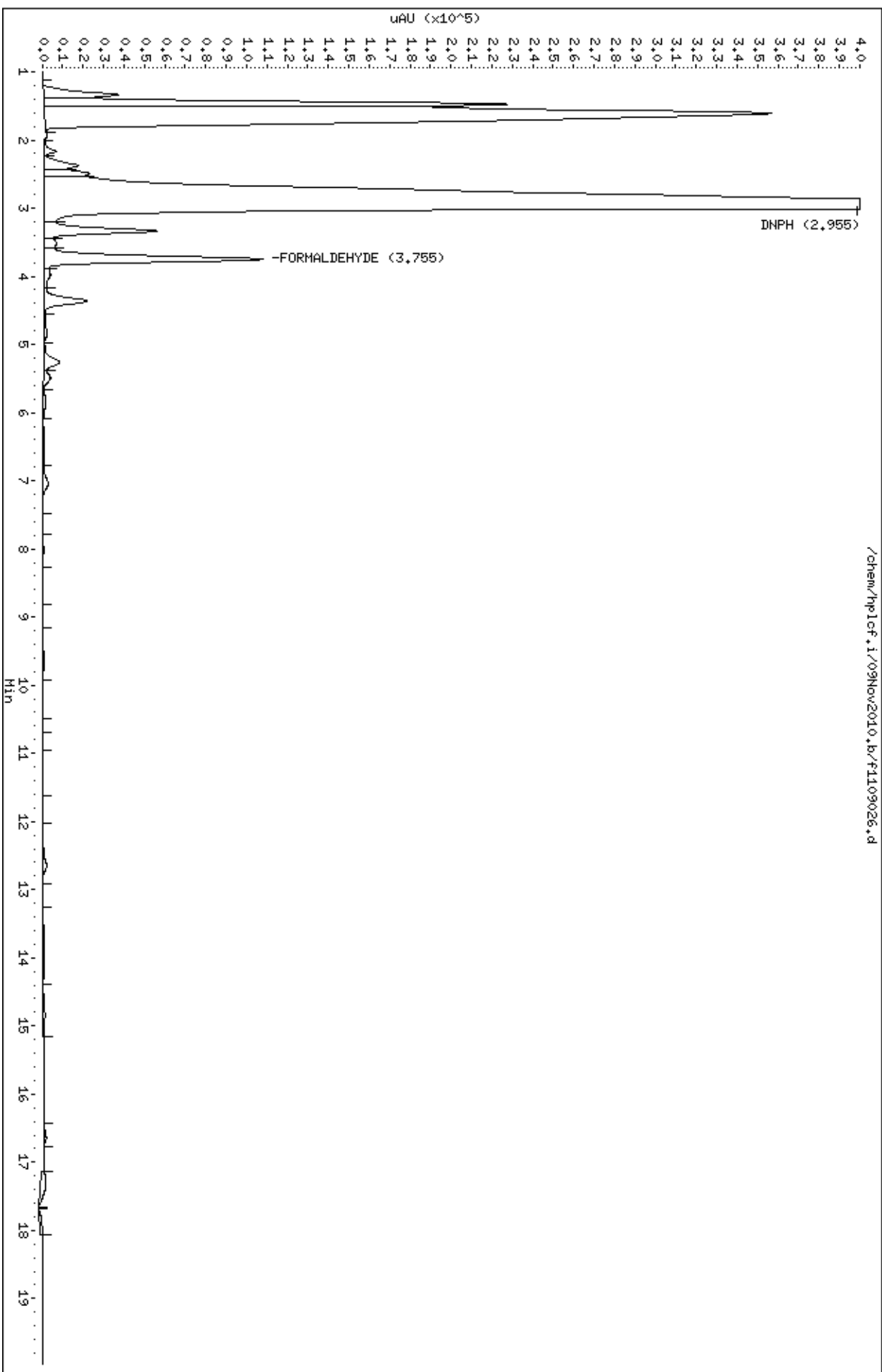
Sample Info: j1011114B-24A;

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



**Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV**

Client Sample ID: 118655

Lab ID#: 1011114B-25A

No Detections Were Found.

Client Sample ID: 118655

Lab ID#: 1011114B-25A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1109021
Dil. Factor: 1.00

Date of Collection: NA
Date of Analysis: 11/9/10 03:35 PM
Date of Extraction: 11/9/10

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	1.0	Not Detected

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109021.d
Lab Smp Id: 1011114B-25A
Inj Date : 09-NOV-2010 15:35
Operator : CRL
Smp Info : ;1011114B-25A;
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf
Cal Date : 14-SEP-2010 18:30
Als bottle: 1
Dil Factor: 1.00000
Integrator: Falcon
Target Version: 3.50
Processing Host: eeyore

Inst ID: hplcf.i

Quant Type: ESTD

Cal File: f0914014.d

Compound Sublist: form.sub

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	2.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	CONCENTRATIONS	
							ON-COLUMN (ng)	FINAL (ug)
1 DNPH	2.925	2.957	-0.032			24039365		
2 FORMALDEHYDE						Compound Not Detected.		

Data File: /chem/hp1cf.i/09Nov2010.b/f1109021.d

Page 1

Date : 09-NOV-2010 15:35

Client ID:

Instrument: hp1cf.i

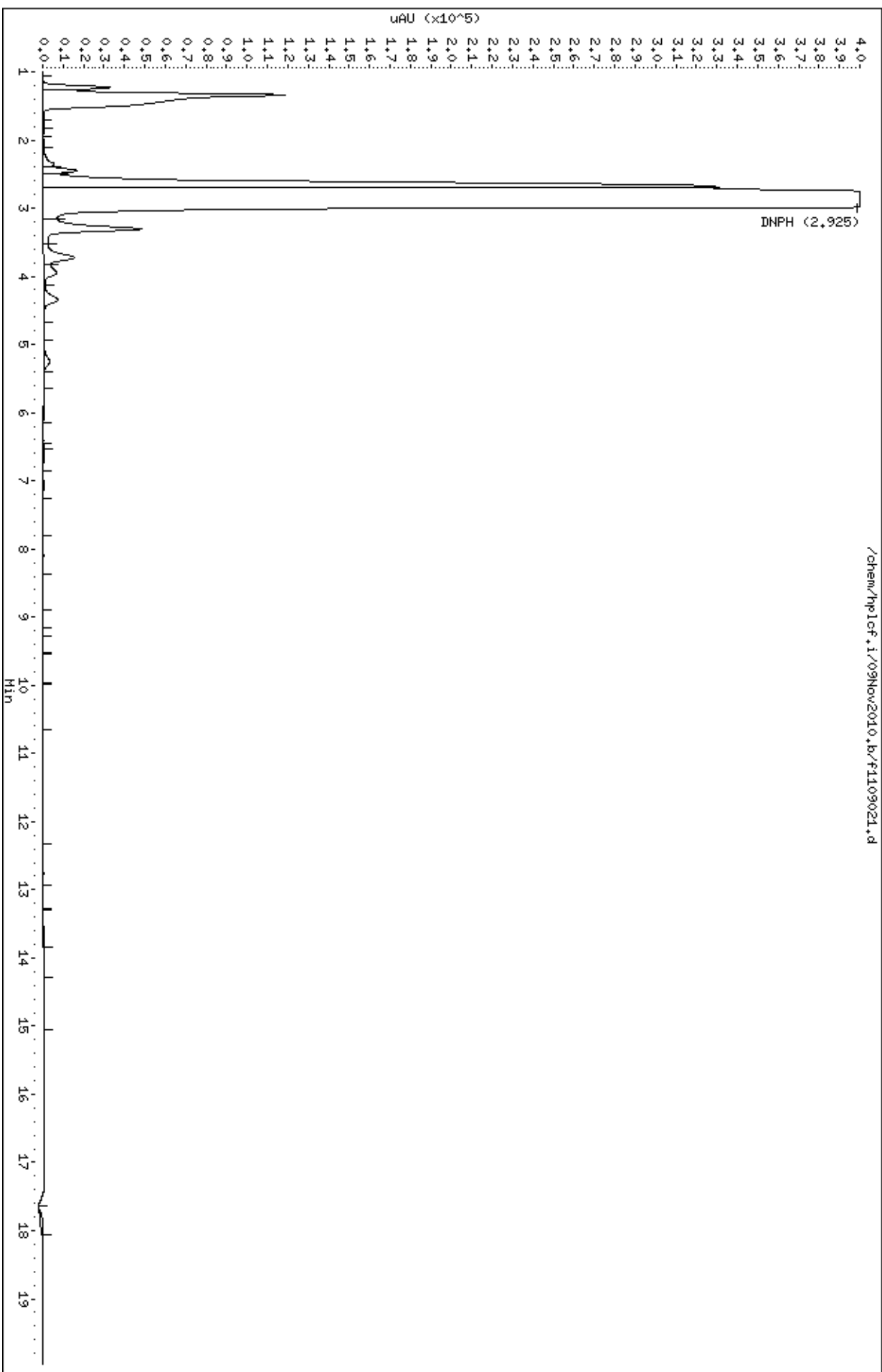
Sample Info: f1011114B-25A;

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



QC Results and Raw Data

Client Sample ID: Lab Blank

Lab ID#: 1011114B-26A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1109010
Dil. Factor: 1.00

Date of Collection: NA
Date of Analysis: 11/9/10 11:45 AM
Date of Extraction: 11/9/10

Compound	Rpt. Limit (ug)	Amount (ug)
Formaldehyde	1.0	Not Detected

Container Type: NA - Not Applicable

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109010.d
Lab Smp Id: 1011114B Client Smp ID: Lab Blank
Inj Date : 09-NOV-2010 11:45
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1011114B;Lab Blank
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: form.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	2.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					CONCENTRATIONS	
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.925	2.957	-0.032	28937954		
2 FORMALDEHYDE				Compound Not Detected.		

Data File: /chem/hp1cf.i/09Nov2010.b/f1109010.d

Page 1

Date : 09-NOV-2010 11:45

Client ID: Lab Blank

Instrument: hp1cf.i

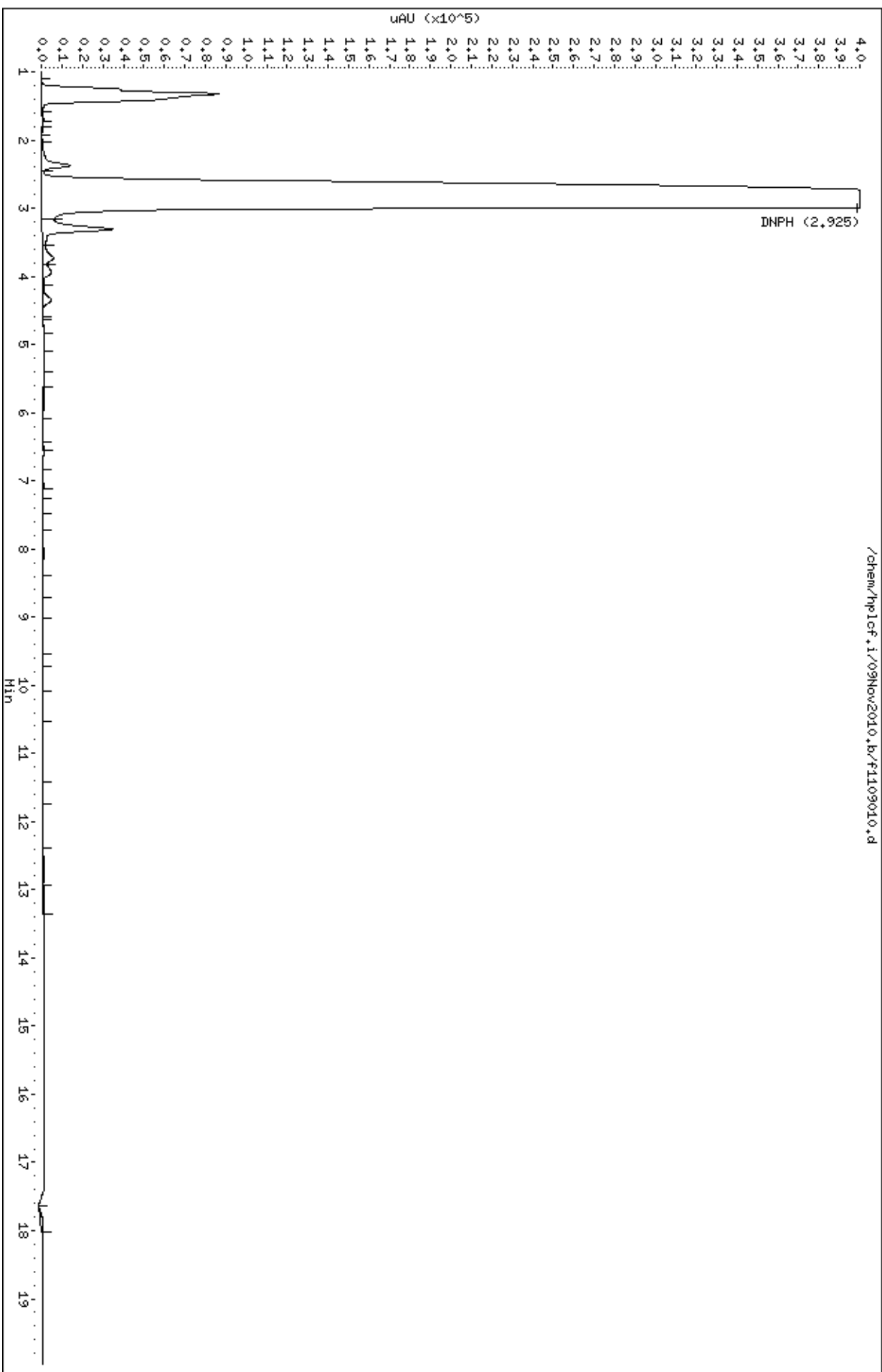
Sample Info: f101114B;Lab Blank

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



SAMPLE RESULTS/SAMPLE RESULTS DUPLICATE

Lab Name: Air Toxics Ltd.

Lab File ID: f1109012.d & f1109011.d

Lab Sample ID: &

Dilution: 1.00 & 1.00

Client Sample ID: LCS & LCSD

Date Analyzed: 11/9/10 & 11/9/10

CAS Number	Compound	Original		Duplicate		RPD	Result Less Than 5X RL
		Amount	Flags	Amount	Flags		
50-00-0	Formaldehyde	117		117		0	

Note: The results appearing in the Amount columns are the raw, unrounded numbers acquired from the instrument.

Air Toxics Ltd.

INITIAL CALIBRATION DATA

Start Cal Date : 14-SEP-2010 16:46
 End Cal Date : 15-SEP-2010 02:52
 Quant Method : ESTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : Falcon
 Method file : /chem/hplcf.i/14Sep2010.b/F10D0914.m
 Cal Date : 15-Sep-2010 09:40 cleaf
 Curve Type : Average

Calibration File Names:

Level 1: /chem/hplcf.i/14Sep2010.b/f0914011.d
 Level 2: /chem/hplcf.i/14Sep2010.b/f0914014.d
 Level 3: /chem/hplcf.i/14Sep2010.b/f0914017.d
 Level 4: /chem/hplcf.i/14Sep2010.b/f0914020.d
 Level 5: /chem/hplcf.i/14Sep2010.b/f0914023.d
 Level 6: /chem/hplcf.i/14Sep2010.b/f0914026.d
 Level 7: /chem/hplcf.i/14Sep2010.b/f0914029.d
 Level 8: /chem/hplcf.i/14Sep2010.b/f0914032.d
 Level 9: /chem/hplcf.i/14Sep2010.b/f0914035.d
 Level 10: /chem/hplcf.i/14Sep2010.b/f0914038.d

Compound	0.05000	0.10000	0.25000	0.50000	1.000	4.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	10.000	20.000	50.000	75.000				
	Level 7	Level 8	Level 9	Level 10				
1 DNPH	+++++	+++++	+++++	+++++	+++++	+++++	+++++	+++++
2 FORMALDEHYDE	123853	105887	102244	90559	105563	112382	109474	8.163
3 ACETALDEHYDE	+++++	84153	81685	84358	82946	88016	85159	3.051
4 ACRYLEIN	+++++	+++++	76745	81228	78019	83736	81132	3.696
5 ACETONE	+++++	+++++	60665	62663	60964	63905	62324	2.766
6 PROPANAL	+++++	+++++	60604	62370	61076	64615	62684	3.112

Air Toxics Ltd.

INITIAL CALIBRATION DATA

Start Cal Date : 14-SEP-2010 16:46
 End Cal Date : 15-SEP-2010 02:52
 Quant Method : ESTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : Falcon
 Method file : /chem/hplcf.i/14Sep2010.b/F10D0914.m
 Cal Date : 15-Sep-2010 09:40 cleaf
 Curve Type : Average

Compound	0.05000 Level 1	0.10000 Level 2	0.25000 Level 3	0.50000 Level 4	1.000 Level 5	4.000 Level 6	RRF	% RSD
	10.000 Level 7	20.000 Level 8	50.000 Level 9	75.000 Level 10				
7 CROTONAL	+++++	+++++	55455	58038	55288	58300		
	54847	56743	57118	59765			56944	2.948
9 BUTYRALS/MEK	+++++	+++++	50991	54916	52929	55652		
	52447	54263	54682	57186			54133	3.608
10 BENZALDEHYDE	+++++	+++++	37551	37827	36952	39328		
	37138	38317	38582	40377			38259	3.289
11 ISO-PENTANAL	+++++	+++++	43787	45109	43981	46732		
	43928	45776	46097	48239			45456	3.516
12 PENTANAL	+++++	+++++	42595	45053	43361	45852		
	43050	44975	45307	47409			44700	3.561
14 O-TOLUALDEHYDE	+++++	+++++	37437	37216	35100	36651		
	34745	36139	36359	38081			36466	3.147
15 M,P-TOLUALDEHYDE	+++++	+++++	31477	31077	29474	31144		
	29314	30236	30483	31864			30634	3.179
16 HEXANAL	+++++	+++++	42072	41339	39329	40825		
	38336	39490	39729	41553			40334	3.839
17 2,5-Dimethylbenzaldehyde	+++++	+++++	32847	31021	28793	29408		
	27544	28236	28401	29711			29495	5.961

Calibration History

Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Start Cal Date: 14-SEP-2010 16:46
End Cal Date : 15-SEP-2010 02:52

Initial Calibration

Injection Date	Sublist	Calibration File
Cal Level: 1 , Cal Amount: 0.05000		
14-SEP-2010 17:27	form	/chem/hplcf.i/14Sep2010.b/f0914011.d
14-SEP-2010 17:07	form	/chem/hplcf.i/14Sep2010.b/f0914010.d
14-SEP-2010 16:46	form	/chem/hplcf.i/14Sep2010.b/f0914009.d

Cal Level: 2 , Cal Amount: 0.10000		
14-SEP-2010 18:30	F+A+M,P	/chem/hplcf.i/14Sep2010.b/f0914014.d
14-SEP-2010 18:09	F+A+M,P	/chem/hplcf.i/14Sep2010.b/f0914013.d
14-SEP-2010 17:48	F+A+M,P	/chem/hplcf.i/14Sep2010.b/f0914012.d

Cal Level: 3 , Cal Amount: 0.25000		
14-SEP-2010 19:33	all	/chem/hplcf.i/14Sep2010.b/f0914017.d
14-SEP-2010 19:12	all	/chem/hplcf.i/14Sep2010.b/f0914016.d
14-SEP-2010 18:51	all	/chem/hplcf.i/14Sep2010.b/f0914015.d

Cal Level: 4 , Cal Amount: 0.50000		
14-SEP-2010 20:35	all	/chem/hplcf.i/14Sep2010.b/f0914020.d
14-SEP-2010 20:15	all	/chem/hplcf.i/14Sep2010.b/f0914019.d
14-SEP-2010 19:54	all	/chem/hplcf.i/14Sep2010.b/f0914018.d

Cal Level: 5 , Cal Amount: 1.00000		
14-SEP-2010 21:38	all	/chem/hplcf.i/14Sep2010.b/f0914023.d
14-SEP-2010 21:17	all	/chem/hplcf.i/14Sep2010.b/f0914022.d
14-SEP-2010 20:56	all	/chem/hplcf.i/14Sep2010.b/f0914021.d

Cal Level: 6 , Cal Amount: 4.00000		
14-SEP-2010 22:41	all	/chem/hplcf.i/14Sep2010.b/f0914026.d
14-SEP-2010 22:20	all	/chem/hplcf.i/14Sep2010.b/f0914025.d

14-SEP-2010 21:59 all	/chem/hplcf.i/14Sep2010.b/f0914024.d	
+-----+		
Cal Level: 7 , Cal Amount: 10.00000		
+=====+		
14-SEP-2010 23:44 all	/chem/hplcf.i/14Sep2010.b/f0914029.d	
14-SEP-2010 23:23 all	/chem/hplcf.i/14Sep2010.b/f0914028.d	
14-SEP-2010 23:02 all	/chem/hplcf.i/14Sep2010.b/f0914027.d	
+-----+		
Cal Level: 8 , Cal Amount: 20.00000		
+=====+		
15-SEP-2010 00:46 all	/chem/hplcf.i/14Sep2010.b/f0914032.d	
15-SEP-2010 00:25 all	/chem/hplcf.i/14Sep2010.b/f0914031.d	
15-SEP-2010 00:04 all	/chem/hplcf.i/14Sep2010.b/f0914030.d	
+-----+		
Cal Level: 9 , Cal Amount: 50.00000		
+=====+		
15-SEP-2010 01:49 all	/chem/hplcf.i/14Sep2010.b/f0914035.d	
15-SEP-2010 01:28 all	/chem/hplcf.i/14Sep2010.b/f0914034.d	
15-SEP-2010 01:07 all	/chem/hplcf.i/14Sep2010.b/f0914033.d	
+-----+		
Cal Level: 10, Cal Amount: 75.00000		
+=====+		
15-SEP-2010 02:52 all	/chem/hplcf.i/14Sep2010.b/f0914038.d	
15-SEP-2010 02:31 all	/chem/hplcf.i/14Sep2010.b/f0914037.d	
15-SEP-2010 02:10 all	/chem/hplcf.i/14Sep2010.b/f0914036.d	
+-----+		

Continuing Calibration

Ccal Level Mode: GLOBAL LEVEL 6

+-----+		
Ccal Level: 6 , Ccal Amount: 4.00		
+=====+		
14-SEP-2010 22:41 all	/chem/hplcf.i/14Sep2010.b/f0914026.d	
+-----+		
Ccal Level: 6 , Ccal Amount: 4.00		
+=====+		
14-SEP-2010 22:20 all	/chem/hplcf.i/14Sep2010.b/f0914025.d	
+-----+		
Ccal Level: 6 , Ccal Amount: 4.00		
+=====+		
14-SEP-2010 21:59 all	/chem/hplcf.i/14Sep2010.b/f0914024.d	
+-----+		

Air Toxics Ltd.

INITIAL CALIBRATION DATA

Start Cal Date : 14-SEP-2010 16:46
 End Cal Date : 15-SEP-2010 02:52
 Quant Method : ESTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : Falcon
 Method file : /chem/hplcf.i/14Sep2010.b/F10D0914.m
 Cal Date : 15-Sep-2010 09:40 cleaf
 Curve Type : Average

Carry over 9/20/10

2nd source : f0914040.d

Calibration File Names:

Level 1: /chem/hplcf.i/14Sep2010.b/f0914011.d
 Level 2: /chem/hplcf.i/14Sep2010.b/f0914014.d
 Level 3: /chem/hplcf.i/14Sep2010.b/f0914017.d
 Level 4: /chem/hplcf.i/14Sep2010.b/f0914020.d
 Level 5: /chem/hplcf.i/14Sep2010.b/f0914023.d
 Level 6: /chem/hplcf.i/14Sep2010.b/f0914026.d
 Level 7: /chem/hplcf.i/14Sep2010.b/f0914029.d
 Level 8: /chem/hplcf.i/14Sep2010.b/f0914032.d
 Level 9: /chem/hplcf.i/14Sep2010.b/f0914035.d
 Level 10: /chem/hplcf.i/14Sep2010.b/f0914038.d

See history

Based on a 20ul injection

9/20/10

Compound	0.05000	0.10000	0.25000	0.50000	1.000	4.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
<i>ug/5ml</i>	10.000	20.000	50.000	75.000				
	Level 7	Level 8	Level 9	Level 10				
1 DNPH	+++++	+++++	+++++	+++++	+++++	+++++		
	+++++	+++++	+++++	+++++			+++++	+++++
2 FORMALDEHYDE	✓123853	✓105887	✓102244	✓90559	✓105563	✓112382		
	✓108702	✓112707	✓114514	✓118333			✓109474	✓8.163
3 ACETALDEHYDE	+++++	84153	81685	84358	82946	88016		
	83072	85795	86254	90148			85159	3.051
4 ACROLEIN	+++++	+++++	76745	81228	78019	83736		
	78948	81614	82596	86166			81132	3.696
5 ACETONE	+++++	+++++	60665	62663	60964	63905		
	60075	62505	62587	65230			62324	2.766
6 PROPANAL	+++++	+++++	60604	62370	61076	64615		
	60743	62795	62980	66291			62684	3.112

Carry over 9/20/10

Air Toxics Ltd.

INITIAL CALIBRATION DATA

Start Cal Date : 14-SEP-2010 16:46
 End Cal Date : 15-SEP-2010 02:52
 Quant Method : ESTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : Falcon
 Method file : /chem/hplcf.i/14Sep2010.b/F10D0914.m
 Cal Date : 15-Sep-2010 09:40 cleaf
 Curve Type : Average

Compound	0.05000	0.10000	0.25000	0.50000	1.000	4.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
	10.000	20.000	50.000	75.000				
	Level 7	Level 8	Level 9	Level 10				
7 CROTONAL	+++++	+++++	55455	58038	55288	58300		
	54847	56743	57118	59765			56944	2.948
9 BUTYRALS/MEK	+++++	+++++	50991	54916	52929	55652		
	52447	54263	54682	57186			54133	3.608
10 BENZALDEHYDE	+++++	+++++	37551	37827	36952	39328		
	37138	38317	38582	40377			38259	3.289
11 ISO-PENTANAL	+++++	+++++	43787	45109	43981	46732		
	43928	45776	46097	48239			45456	3.516
12 PENTANAL	+++++	+++++	42595	45053	43361	45852		
	43050	44975	45307	47409			44700	3.561
14 O-TOLUALDEHYDE	+++++	+++++	37437	37216	35100	36651		
	34745	36139	36359	38081			36466	3.147
15 M, P-TOLUALDEHYDE	+++++	+++++	31477	31077	29474	31144		
	29314	30236	30483	31864			30634	3.179
16 HEXANAL	+++++	+++++	42072	41339	39329	40825		
	38336	39490	39729	41553			40334	3.839
17 2,5-Dimethylbenzaldehyde	+++++	+++++	32847	31021	28793	29408		
	27544	28236	28401	29711			29495	5.961

Method: All Methods (DNPH)Pressure: 87 bar

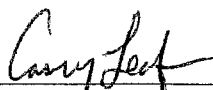
U s e	File #	Sample Name/Client	DF	Loaded by Init.	Date	Time	Reviewed by Init.	Comments
✓	0914008	ACN WASH	1.00	CRL	9/14/10	1625	CRL	equilibrating system
✓	9	1838-34-0.05				1625 1646		level 1
✓	10	↓				1707		↓
✓	11	↓				1727		↓
✓	12	1838-34-0.1				1748		level 2
✓	13	↓				1809		↓
✓	14	↓				1830		↓
✓	15	1838-34-0.25				1851		level 3
✓	16	↓				1912		↓
✓	17	↓				1933		↓
✓	18	1838-34-0.50				1954		level 4
✓	19	↓				2015		↓
✓	20	↓				2035		↓
✓	21	1838-34-1.0				2056		level 5
✓	22	↓				2117		↓
✓	23	↓				2138		↓
✓	24	1838-32-4.0				239		level 6
✓	25	↓				2220		↓
✓	26	↓				2241		↓
✓	27	1838-34-10.0				2302		level 7
✓	28	↓				2323		↓
✓	29	↓				2344		↓
✓	30	1838-34-20.0			9/15/10	0004		level 8
✓	31	↓				0025		↓
✓	32	↓				0046		↓
✓	33	1838-34-50.0				0107		level 9
✓	34	↓				0128		↓
✓	35	↓				0147		↓

Calculation Check:File ID: 0914024.d Compound: Cotanol Initials: CRLSample Amt = $\frac{\text{Area Counts Sample}}{\text{RF}} \times \text{DF} = \left(\frac{232736}{56944} \right) \times (1.00) =$

4.09

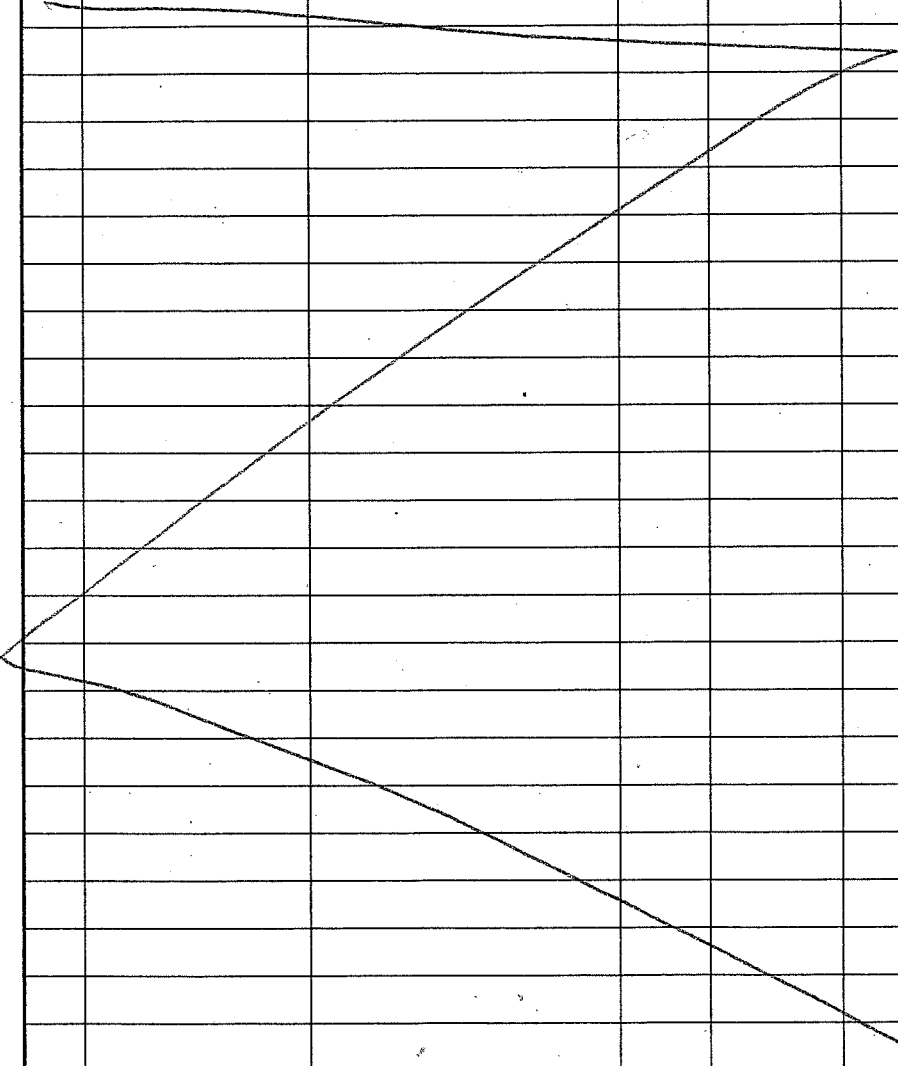
Reported Result:

4.00


 Signed


 Date

Method: All DPH methodsPressure: 87 bar

U s e	File #	Sample Name/Client	DF	Loaded by Init.	Date	Time	Reviewed by Init.	Comments
✓	914036	1838-34-75.0	1.00	CR	9/15/10	0210	CR	level 10
✓	37	↓	↓	↓	↓	0231	↓	↓
✓	38	↓	↓	↓	↓	0252	↓	↓
✓	39	ACU WASH	↓	↓	↓	0312	↓	↓
✓	40	1838-30-4.0 LCS	↓	↓	↓	0333	↓	LCS
								

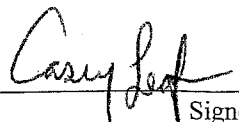
Calculation Check:

File ID: _____ Compound: _____

Initials: _____

Sample Amt = $\frac{\text{Area Counts Sample}}{\text{RF}} \times \text{DF} = \left(\frac{\quad}{\quad} \right) \times \left(\quad \right) =$

Reported Result: _____



Signed

9/15/10

Date

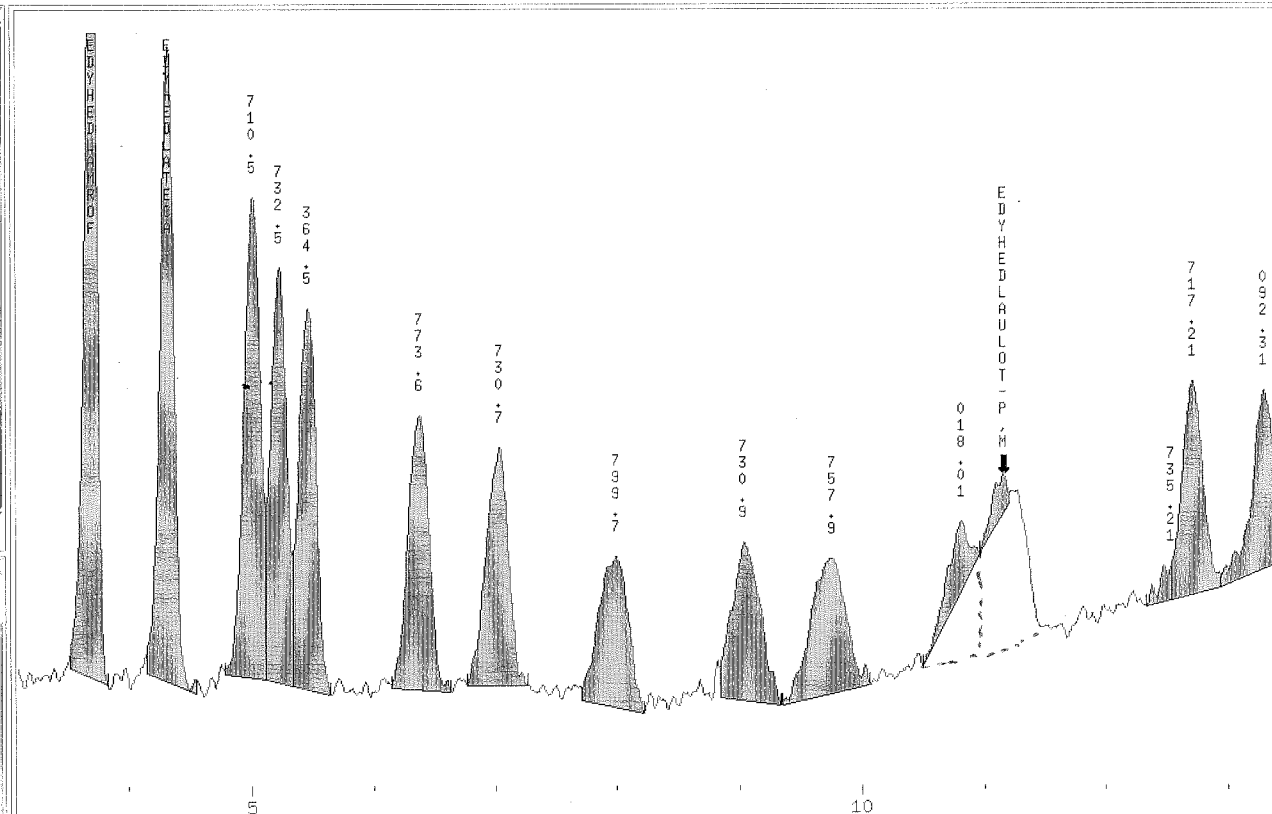
Revised 07/08

File Record Plot Display Process Help

Sample: 1838-34-0.1 Type: CALIB_2 Inj.Date: 14-SEP-2010 18:30 Col: 1 C-18 REVERSE PHASE

1 DNPH
+ 2 FORMALDEHYDE
+ 3 ACETALDEHYDE
+ 15 M,P-TOLUALDEHYDE

F0914011.d
F0914012.d
F0914013.d
F0914014.d
F0914015.d
F0914016.d
F0914017.d
F0914018.d
F0914019.d
F0914020.d
F0914021.d
F0914022.d
F0914023.d
F0914027.d
F0914028.d
F0914029.d
F0914030.d
F0914031.d
F0914032.d
F0914033.d
F0914034.d



Hit# RT(min) Response Amount Conc Ratio Flags Report:

1 11.165 15167 0.490 0.490 100

- Mark M,P-TOLUALDEHYDE Undetected.

Before: Assign baseline and split peak

09/15/10

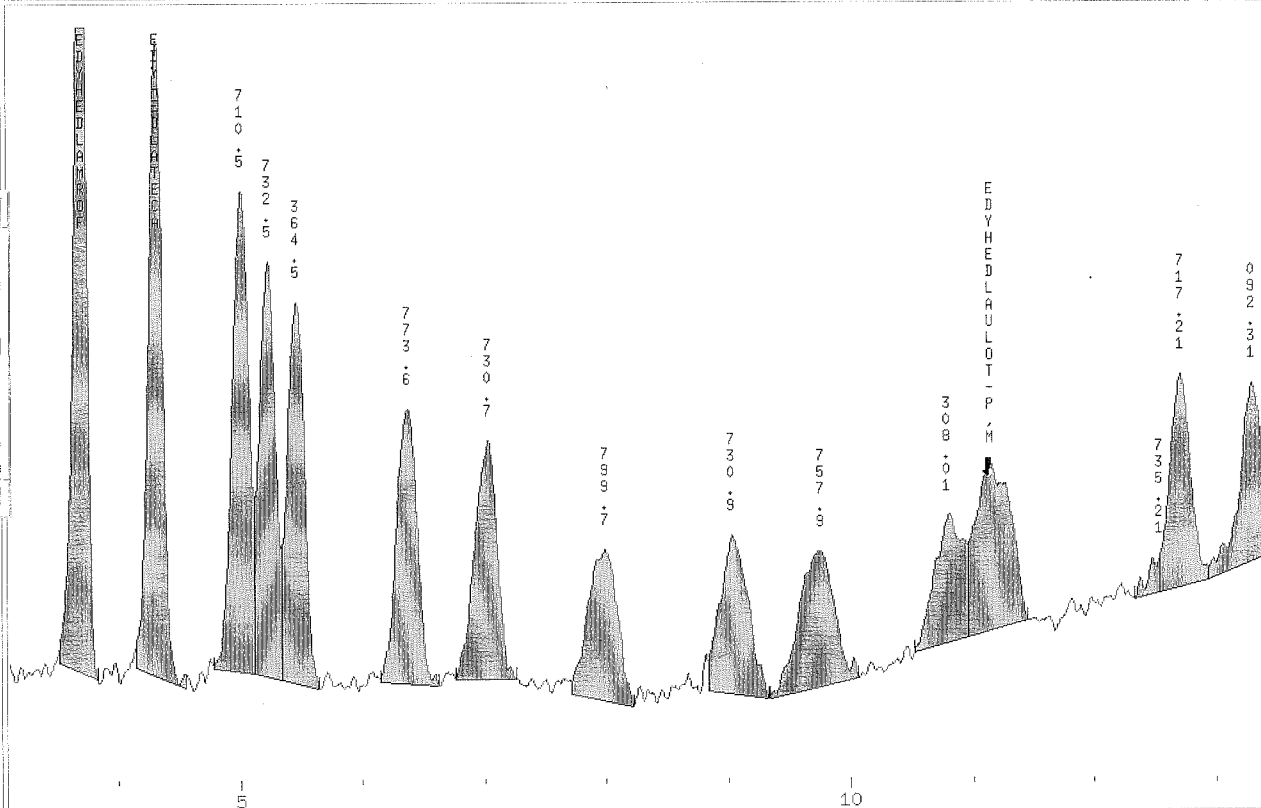
Sample: 1838-34-0.1 Type: CALIB_2 Inj.Date: 14-SEP-2010 18:30 Col: 1 C-18 REVERSE PHASE

1 DNPH
+ 2 FORMALDEHYDE
+ 3 ACETALDEHYDE
+ 15 M,P-TOLUALDEHYDE

Compound: M,P-TOLUAL
Update Method Expected
Time New Time
11.157

Done Apply Una

F0914011.d
F0914012.d
F0914013.d
F0914014.d
F0914015.d
F0914016.d
F0914017.d
F0914018.d
F0914019.d
F0914020.d
F0914021.d
F0914022.d
F0914023.d
F0914027.d
F0914028.d
F0914029.d
F0914030.d
F0914031.d
F0914032.d
F0914033.d
F0914034.d



Hit# RT(min) Response Amount Conc Ratio Flags Report:

1 11.157 5452 0.178 0.178 100 M
- Mark M,P-TOLUALDEHYDE Undetected.

After

Correct Baseline	✓
Split Peak	✓
Merge Peak	N/A
Zoom In	✓
Change Parameter	N/A
System Peak Subtraction	✓
Peak Misidentified	✓
Corrected Peak Integration	✓

09/15/10

09/15/10

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914040.d
Lab Smp Id: 1838-36-4.0 Client Smp ID: LCS
Inj Date : 15-SEP-2010 03:33
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-36-4.0;LCS
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:54 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1 QC Sample: LCS
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					CONCENTRATIONS	
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (UG)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	Compound Not Detected.					
2 FORMALDEHYDE	3.741	3.703	0.038	459327	4.19575	4.20
3 ACETALDEHYDE	4.375	4.323	0.052	346792	4.07232	4.07
4 ACROLEIN	5.095	5.093	0.002	321649	3.96455	3.96
5 ACETONE	5.301	5.300	0.001	259469	4.16321	4.16
6 PROPANAL	5.542	5.546	-0.004	253071	4.03724	4.04
7 CROTONAL	6.488	6.480	0.008	235642	4.13814	4.14
9 BUTYRALS/MEK	7.141	7.146	-0.005	218369	4.03393	4.03
10 BENZALDEHYDE	8.175	8.166	0.009	157886	4.12678	4.13
11 ISO-PENTANAL	9.281	9.260	0.021	183154	4.02926	4.03
12 PENTANAL	9.955	9.946	0.009	174549	3.90489	3.90
14 O-TOLUALDEHYDE	11.035	11.026	0.009	143736	3.94165	3.94
15 M,P-TOLUALDEHYDE	11.341	11.157	0.184	251412	8.20703	8.21
16 HEXANAL	12.861	12.853	0.008	157763	3.91141	3.91
17 2,5-Dimethylbenzaldehyde	13.455	13.453	0.002	115179	3.90508	3.90

Air Toxics Ltd.

RECOVERY REPORT

Client Name: Client SDG: 14Sep2010
Sample Matrix: LIQUID Fraction: OTHER
Lab Smp Id: 1838-36-4.0 Client Smp ID: LCS
Level: MED Operator: CRL
Data Type: LC DATA SampleType: LCS
SpikeList File: 4ug2nd.spk Quant Type: ESTD
Sublist File: all.sub
Method File: /chem/hplcf.i/14Sep2010.b/F10D0914.m
Misc Info:

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	4.00	4.20	104.89	85-115
3 ACETALDEHYDE	4.00	4.07	101.81	85-115
4 ACROLEIN	4.00	3.96	99.11	85-115
5 ACETONE	4.00	4.16	104.08	85-115
6 PROPANAL	4.00	4.04	100.93	85-115
7 CROTONAL	4.00	4.14	103.45	85-115
9 BUTYRALS/MEK	4.00	4.03	100.85	85-115
10 BENZALDEHYDE	4.00	4.13	103.17	85-115
11 ISO-PENTANAL	4.00	4.03	100.73	85-115
12 PENTANAL	4.00	3.90	97.62	85-115
14 O-TOLUALDEHYDE	4.00	3.94	98.54	85-115
15 M,P-TOLUALDEHYDE	8.00	8.21	102.59	85-115
16 HEXANAL	4.00	3.91	97.79	85-115
17 2,5-Dimethylbenzal	4.00	3.90	97.63	85-115

Data File: /chem/hp1cf.i/14Sep2010.b/f0914040.d

Date : 15-SEP-2010 03:33

Client ID: LCS

Sample Info: f1838-36-4.0;LCS

Purge Volume: 5.0

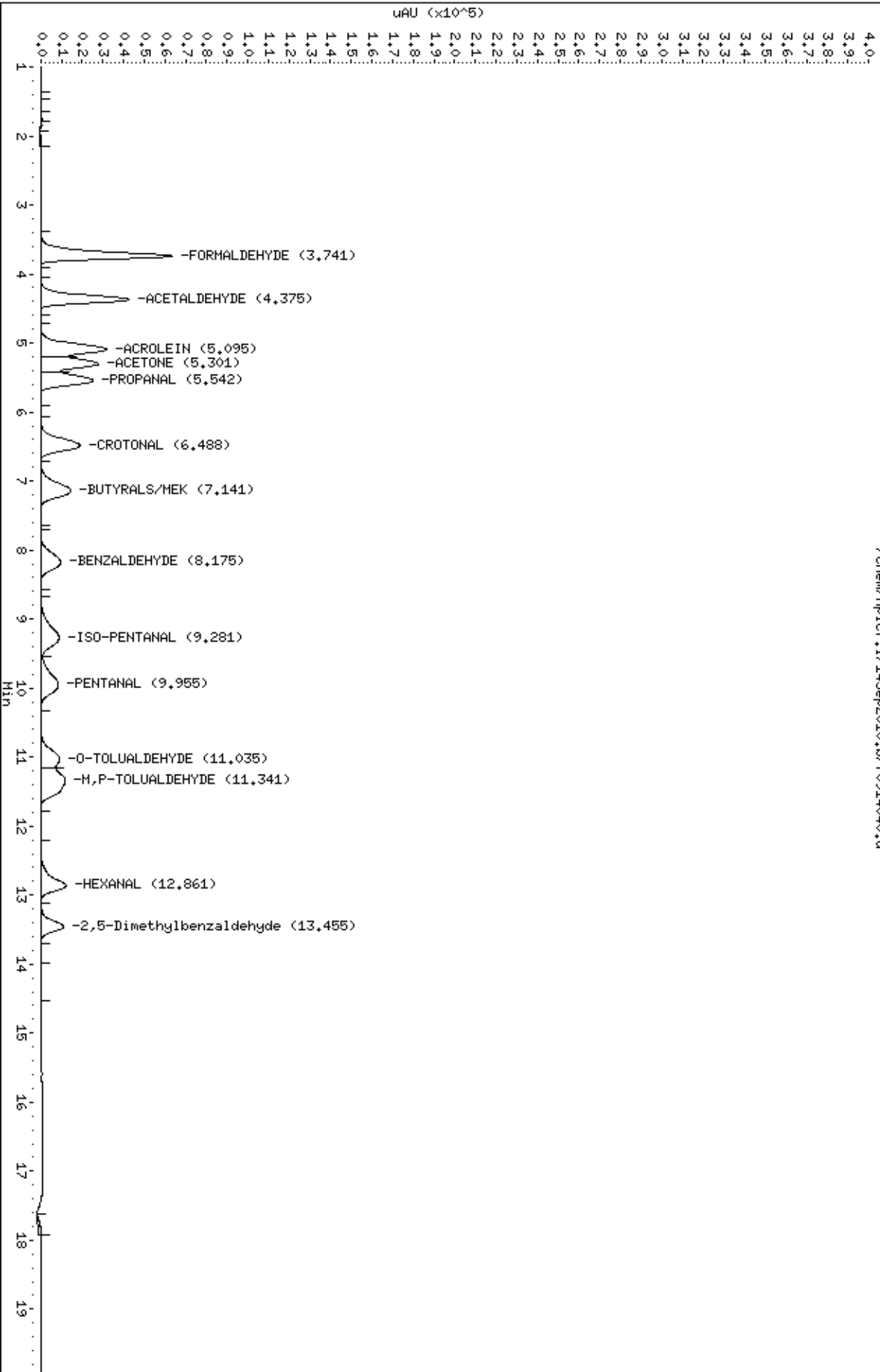
Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00

/chem/hp1cf.i/14Sep2010.b/f0914040.d



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914011.d
Lab Smp Id: 1838-34-0.05 Client Smp ID: 1
Inj Date : 14-SEP-2010 17:27
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-0.05;1
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:39 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 17:27 Cal File: f0914011.d
Als bottle: 1 Calibration Sample, Level: 1
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: form.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH				Compound Not Detected.		
2 FORMALDEHYDE	3.680	3.680	0.000	6159	0.05000	0.0521

Data File: /chem/hplcf.i/14Sep2010.b/f0914011.d

Page 1

Date : 14-SEP-2010 17:27

Client ID: 1

Instrument: hplcf.i

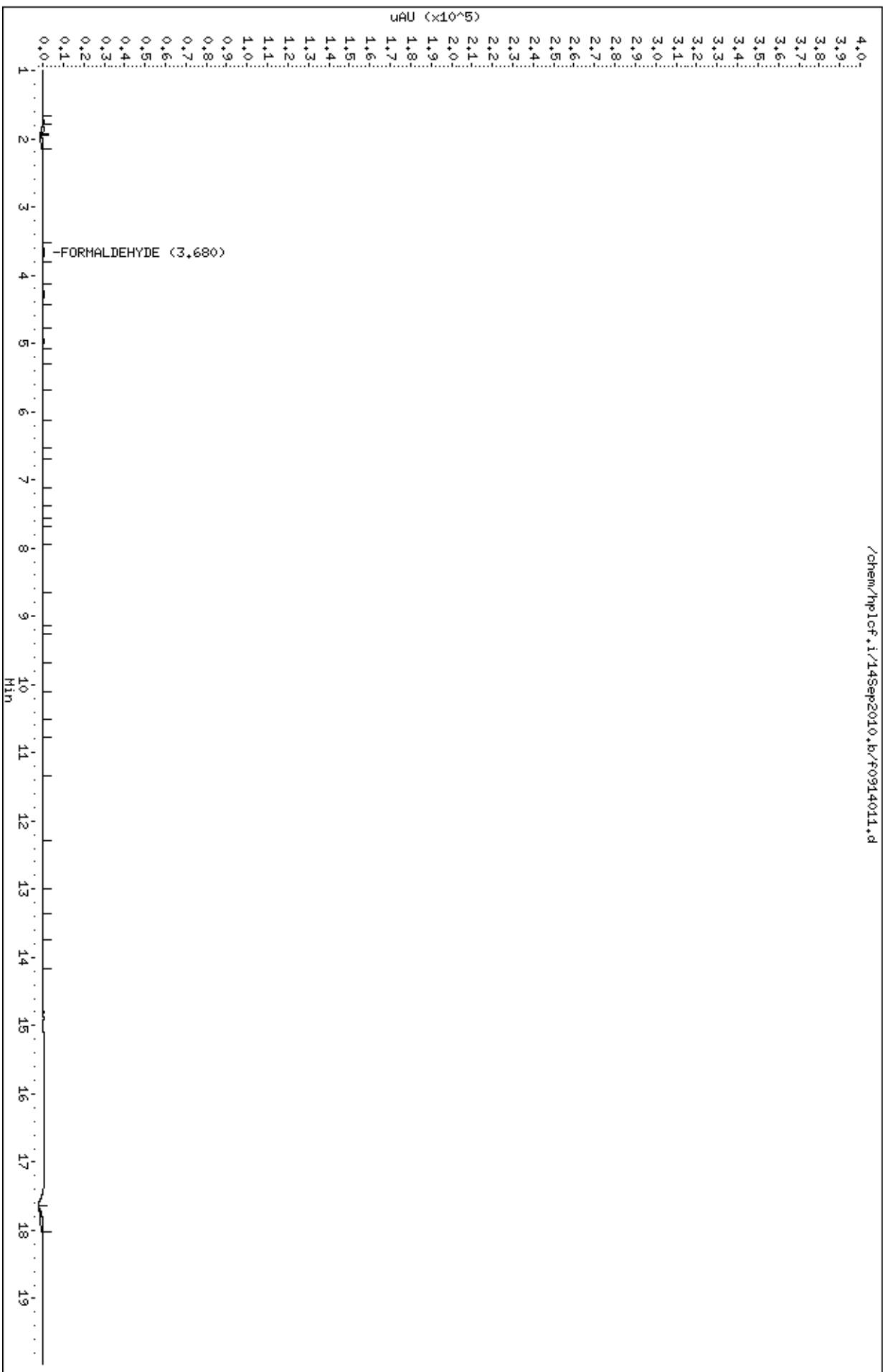
Sample Info: f1838-34-0.05;1

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914010.d
Lab Smp Id: 1838-34-0.05 Client Smp ID: 1
Inj Date : 14-SEP-2010 17:07
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-0.05;1
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:39 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 17:07 Cal File: f0914010.d
Als bottle: 1 Calibration Sample, Level: 1
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: form.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH				Compound Not Detected.		
2 FORMALDEHYDE	3.673	3.673	0.000	6217	0.05000	0.0531

Data File: /chem/hplcf.i/14Sep2010.b/f0914010.d

Page 1

Date : 14-SEP-2010 17:07

Client ID: 1

Instrument: hplcf.i

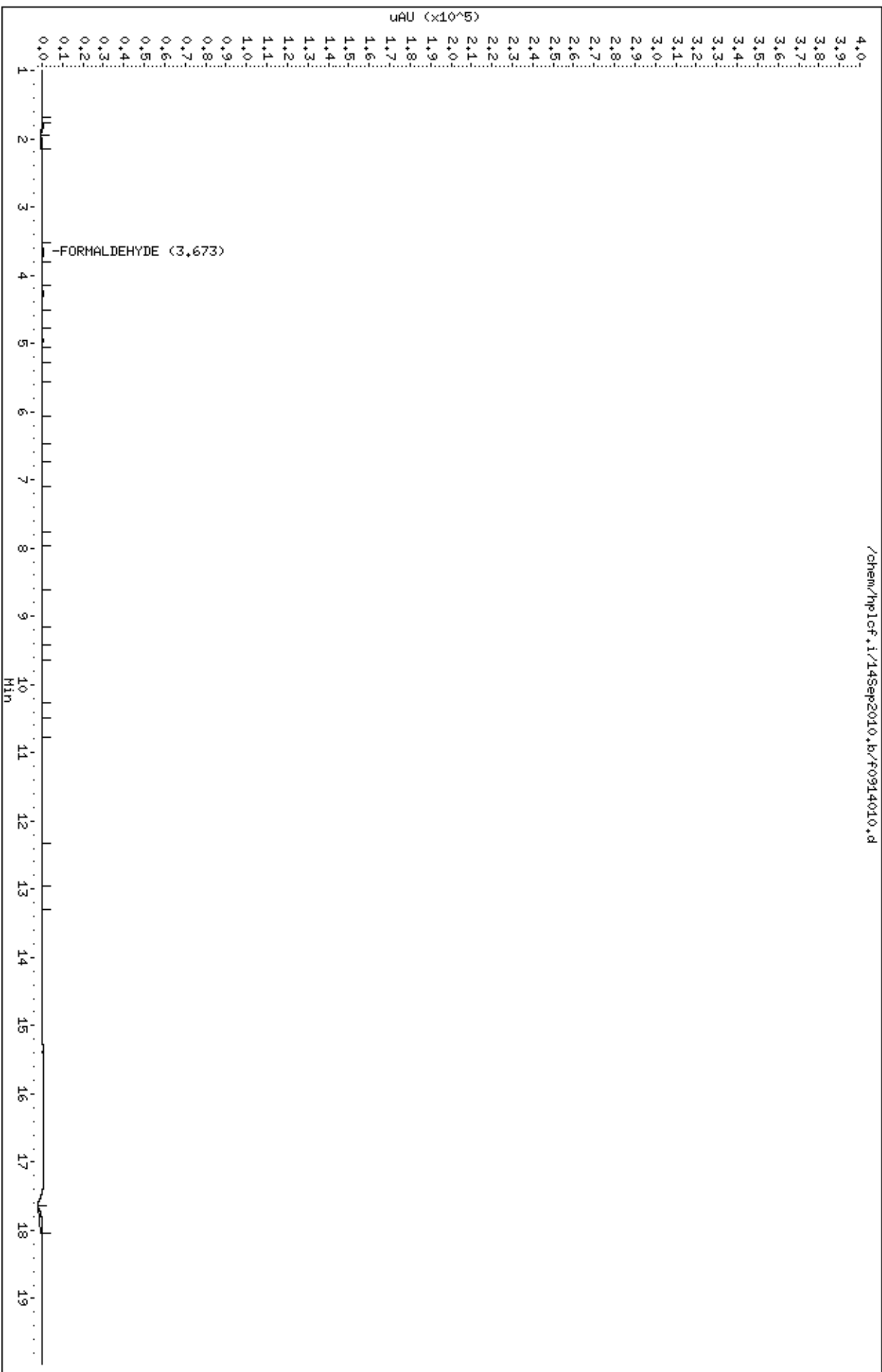
Sample Info: #1838-34-0.05;1

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914009.d
Lab Smp Id: 1838-34-0.05 Client Smp ID: 1
Inj Date : 14-SEP-2010 16:46
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-0.05;1
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:39 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 16:46 Cal File: f0914009.d
Als bottle: 1 Calibration Sample, Level: 1
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: form.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH				Compound Not Detected.		
2 FORMALDEHYDE	3.669	3.669	0.000	6202	0.05000	0.0538

Data File: /chem/hp1cf.i/14Sep2010.b/f0914009.d

Page 1

Date : 14-SEP-2010 16:46

Client ID: 1

Instrument: hp1cf.i

Sample Info: #1838-34-0.05;1

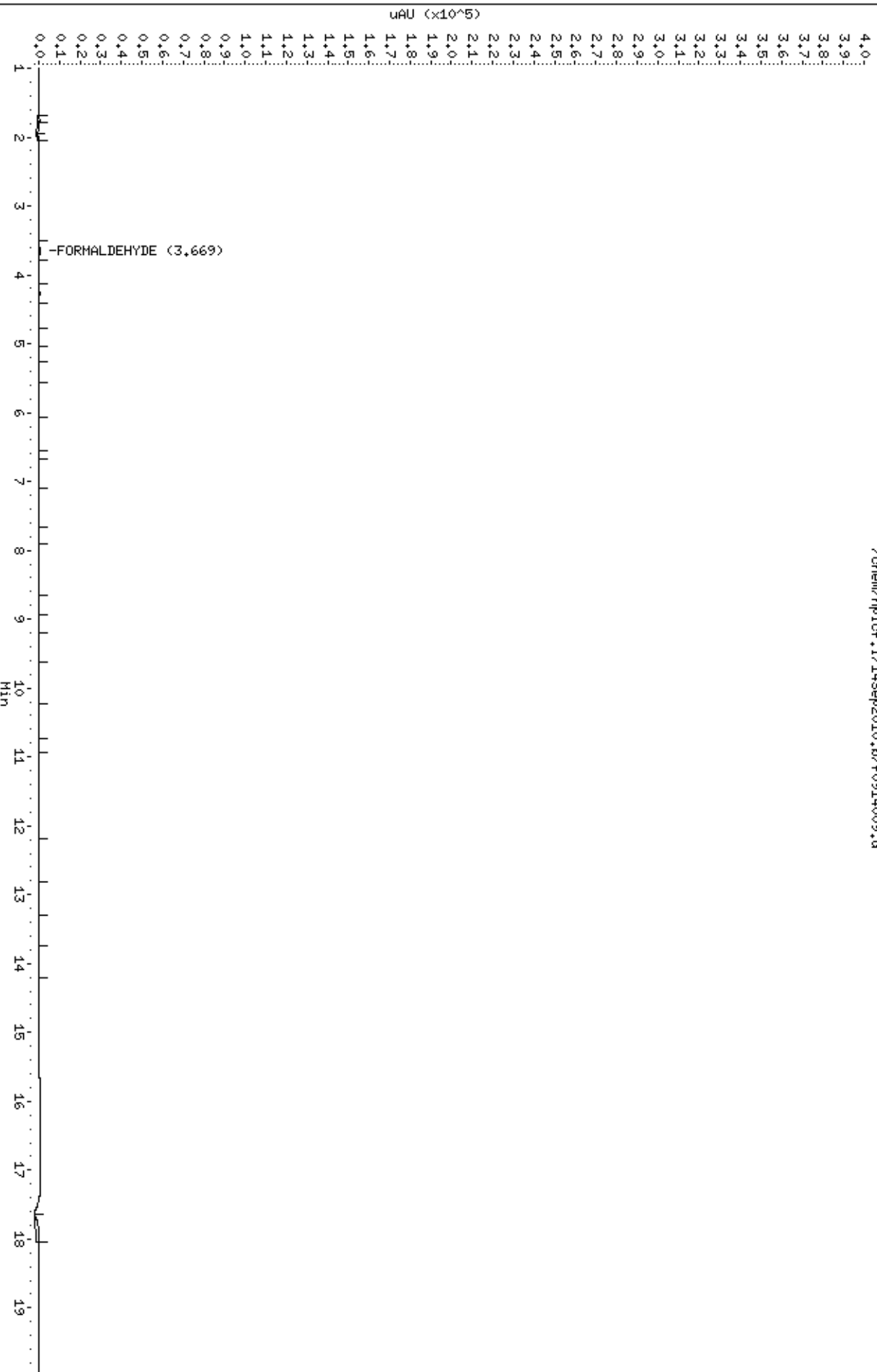
Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00

/chem/hp1cf.i/14Sep2010.b/f0914009.d



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914014.d
Lab Smp Id: 1838-34-0.1 Client Smp ID: 2
Inj Date : 14-SEP-2010 18:30
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-0.1;2
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1 Calibration Sample, Level: 2
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: F+A+M,P.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	AMOUNTS	
							CAL - AMT	ON - COL
							(UG)	(UG)
=====	==	=====	=====			=====	=====	
1 DNPH	Compound Not Detected.							
2 FORMALDEHYDE	3.703	3.740	-0.037		10562	0.10000	0.0965	
3 ACETALDEHYDE	4.323	4.373	-0.050		8338	0.10000	0.0979	
15 M,P-TOLUALDEHYDE	11.157	11.333	-0.176		5452	0.20000	0.178(M)	

QC Flag Legend

M - Compound response manually integrated.

Data File: /chem/hplcf.i/14Sep2010.b/f0914014.d

Date : 14-SEP-2010 18:30

Client ID: 2

Sample Info: f1838-34-0.1;2

Purge Volume: 5.0

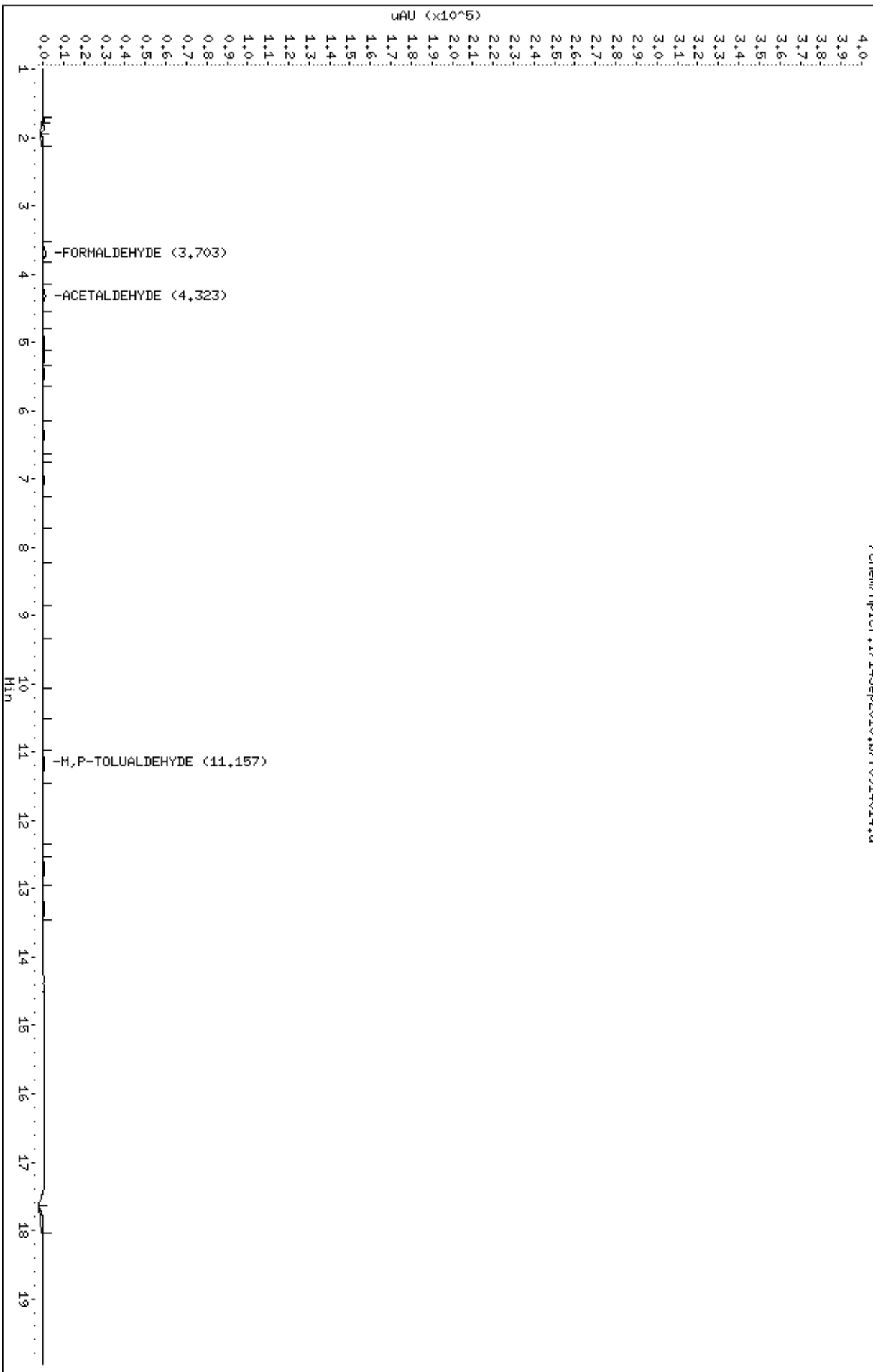
Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: CRL

Column diameter: 2.00

/chem/hplcf.i/14Sep2010.b/f0914014.d



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914013.d
Lab Smp Id: 1838-34-0.1 Client Smp ID: 2
Inj Date : 14-SEP-2010 18:09
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-0.1;2
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:39 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:09 Cal File: f0914013.d
Als bottle: 1 Calibration Sample, Level: 2
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: F+A+M,P.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH	Compound Not Detected.					
2 FORMALDEHYDE	3.704	3.704	0.000	10430	0.10000	0.0906
3 ACETALDEHYDE	4.324	4.324	0.000	8491	0.10000	0.0980
15 M,P-TOLUALDEHYDE	11.124	11.124	0.000	7199	0.20000	0.231

Data File: /chem/hp1cf.i/14Sep2010.b/f0914013.d

Page 1

Date : 14-SEP-2010 18:09

Client ID: 2

Instrument: hp1cf.i

Sample Info: f1838-34-0.1;2

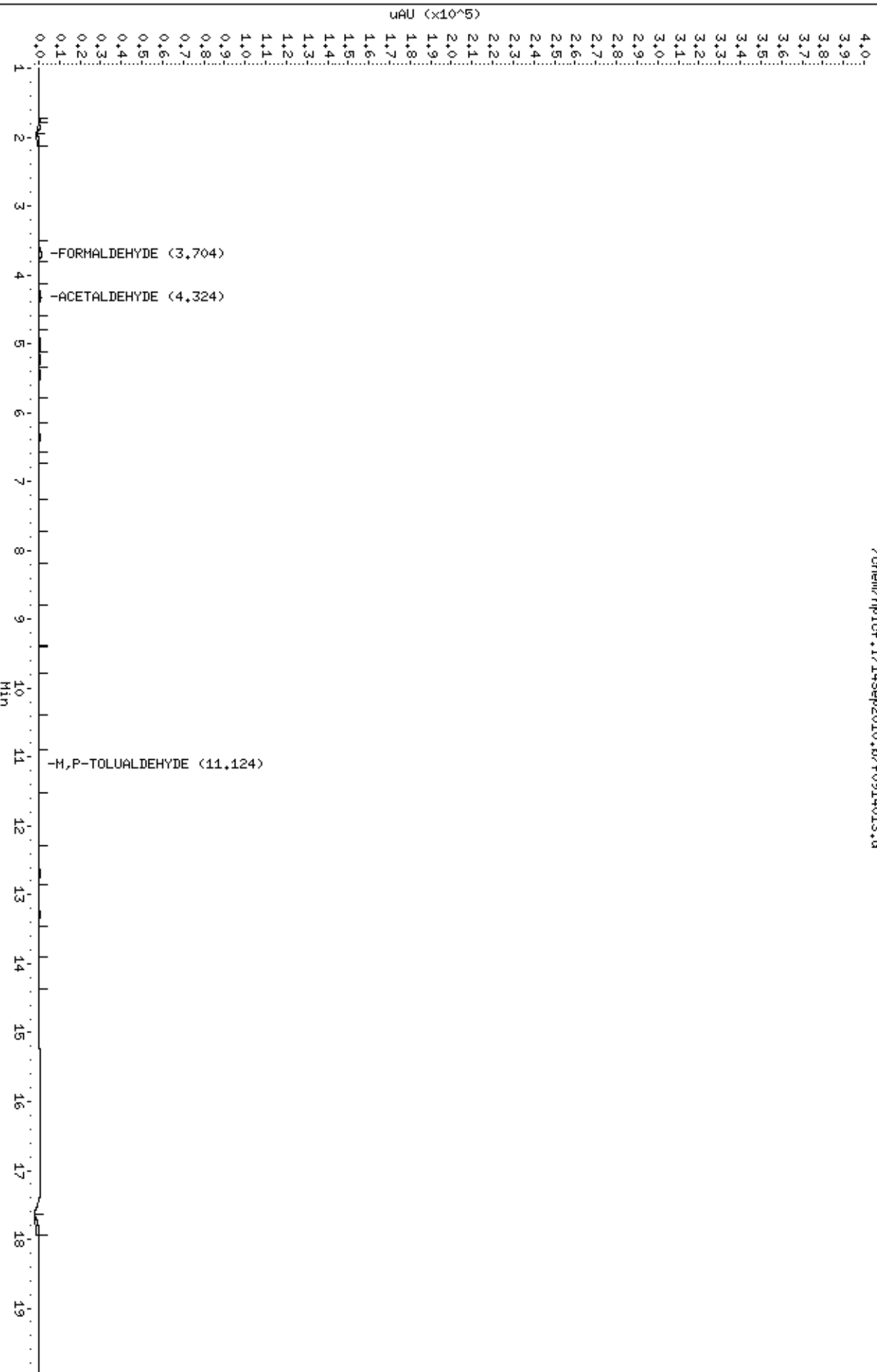
Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00

/chem/hp1cf.i/14Sep2010.b/f0914013.d



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914012.d
Lab Smp Id: 1838-34-0.1 Client Smp ID: 2
Inj Date : 14-SEP-2010 17:48
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-0.1;2
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:39 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 17:48 Cal File: f0914012.d
Als bottle: 1 Calibration Sample, Level: 2
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: F+A+M,P.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH	Compound Not Detected.					
2 FORMALDEHYDE	3.688	3.688	0.000	10774	0.10000	0.0924
3 ACETALDEHYDE	4.301	4.301	0.000	8417	0.10000	0.0967
15 M,P-TOLUALDEHYDE	11.094	11.094	0.000	7091	0.20000	0.228

Data File: /chem/hp1cf.i/14Sep2010.b/f0914012.d

Page 1

Date : 14-SEP-2010 17:48

Client ID: 2

Instrument: hp1cf.i

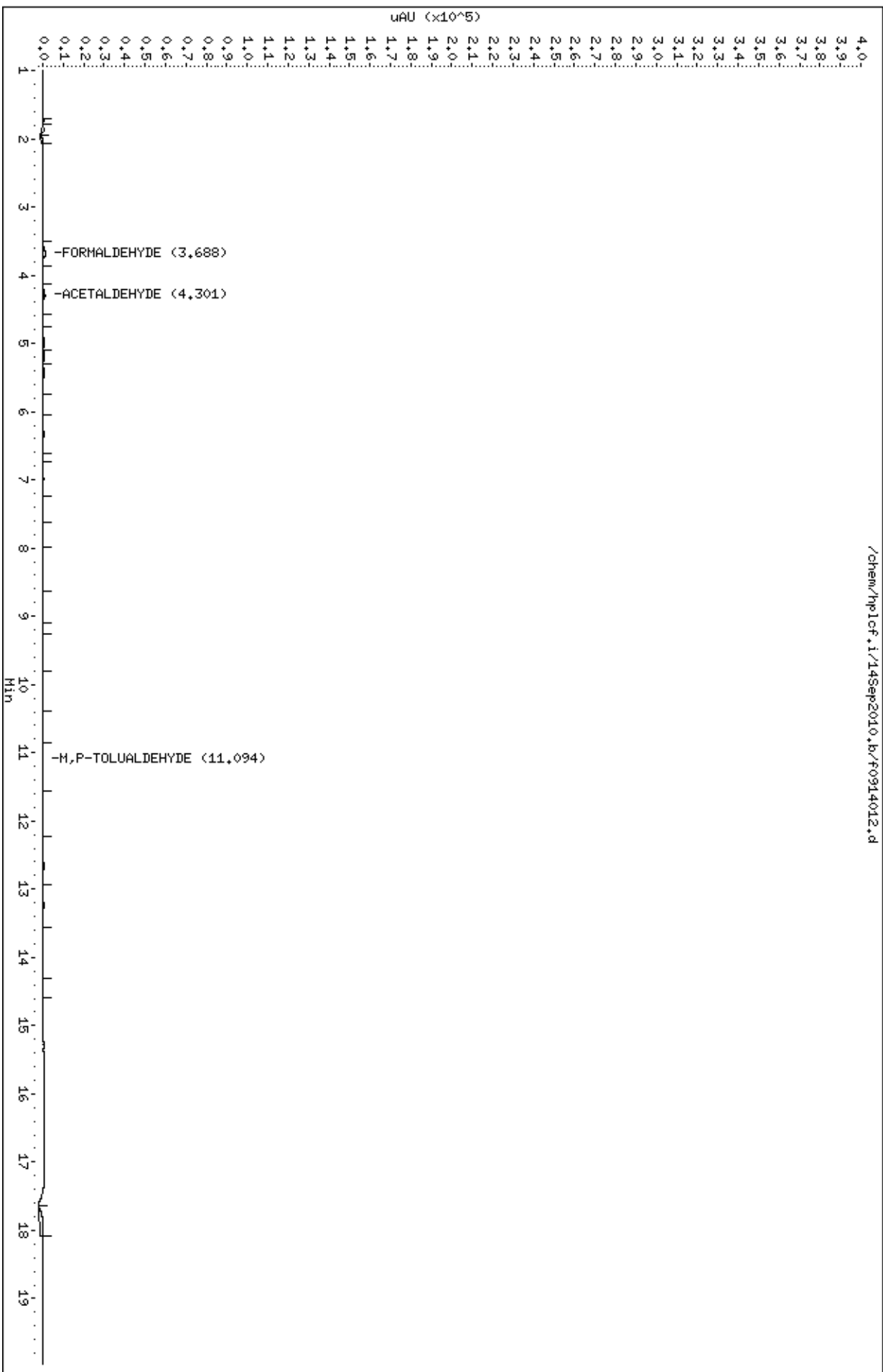
Sample Info: f1838-34-0.1;2

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914017.d
Lab Smp Id: 1838-34-0.25 Client Smp ID: 3
Inj Date : 14-SEP-2010 19:33
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-0.25;3
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:39 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 19:33 Cal File: f0914017.d
Als bottle: 1 Calibration Sample, Level: 3
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH	Compound Not Detected.					
2 FORMALDEHYDE	3.721	3.721	0.000	25225	0.25000	0.227
3 ACETALDEHYDE	4.341	4.341	0.000	20405	0.25000	0.241
4 ACRROLEIN	5.055	5.055	0.000	19200	0.25000	0.239
5 ACETONE	5.261	5.261	0.000	15302	0.25000	0.246
6 PROPANAL	5.508	5.508	0.000	15417	0.25000	0.246
7 CROTONAL	6.428	6.428	0.000	14175	0.25000	0.249
9 BUTYRALS/MEK	7.095	7.095	0.000	12963	0.25000	0.243
10 BENZALDEHYDE	8.108	8.108	0.000	9855	0.25000	0.256
11 ISO-PENTANAL	9.188	9.188	0.000	11199	0.25000	0.247
12 PENTANAL	9.868	9.868	0.000	10971	0.25000	0.248
14 O-TOLUALDEHYDE	10.908	10.908	0.000	9254	0.25000	0.250
15 M,P-TOLUALDEHYDE	11.241	11.241	0.000	15634	0.50000	0.499
16 HEXANAL	12.801	12.801	0.000	10836	0.25000	0.261
17 2,5-Dimethylbenzaldehyde	13.388	13.388	0.000	8638	0.25000	0.278

Data File: /chem/hp1cf.i/14Sep2010.b/f0914017.d

Date : 14-SEP-2010 19:33

Client ID: 3

Sample Info: f1838-34-0.25;3

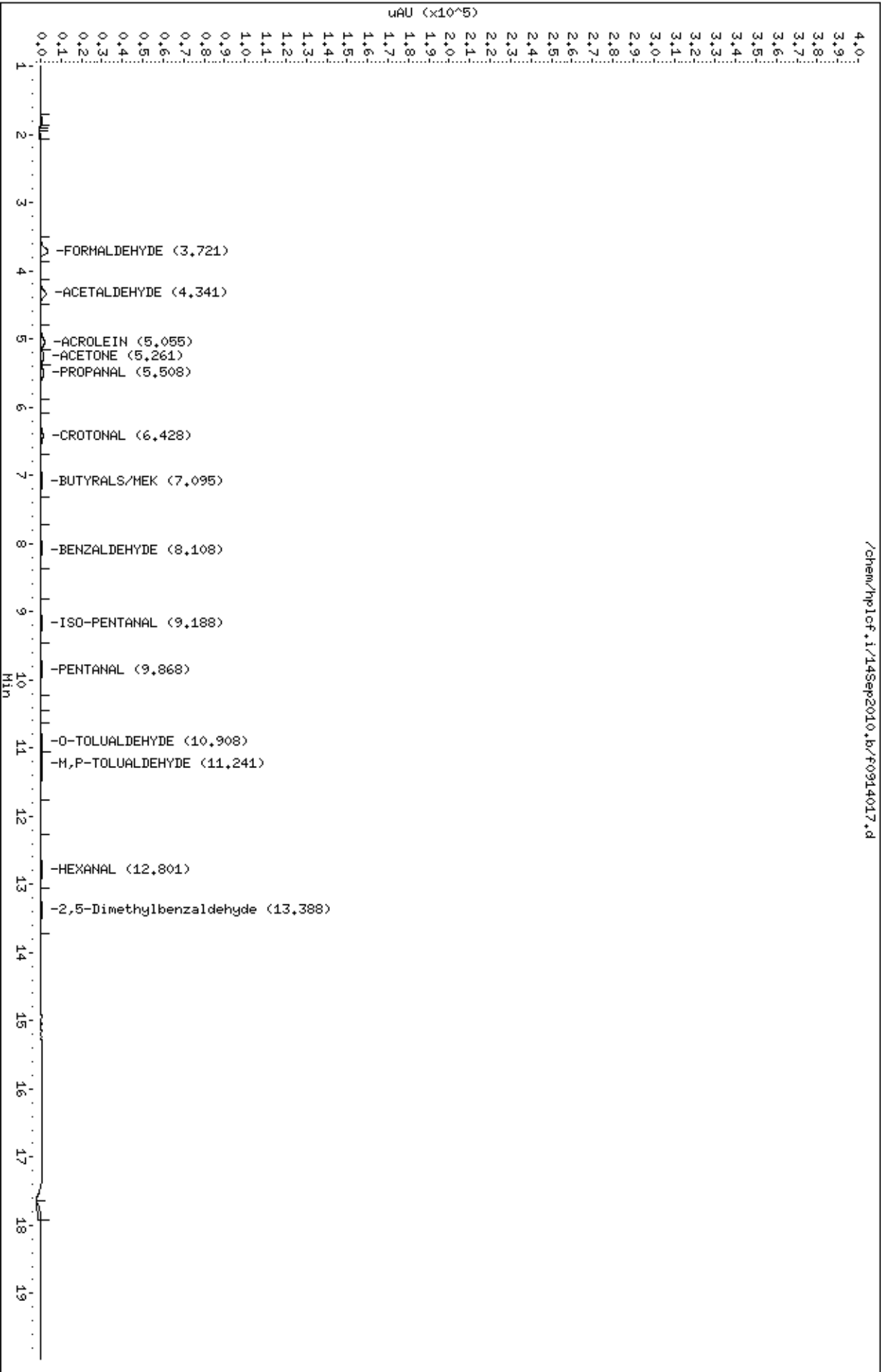
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914016.d
Lab Smp Id: 1838-34-0.25 Client Smp ID: 3
Inj Date : 14-SEP-2010 19:12
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-0.25;3
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:39 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 19:12 Cal File: f0914016.d
Als bottle: 1 Calibration Sample, Level: 3
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH	Compound Not Detected.					
2 FORMALDEHYDE	3.715	3.715	0.000	25598	0.25000	0.228
3 ACETALDEHYDE	4.335	4.335	0.000	20381	0.25000	0.240
4 ACRROLEIN	5.041	5.041	0.000	19129	0.25000	0.236
5 ACETONE	5.255	5.255	0.000	14931	0.25000	0.239
6 PROPANAL	5.495	5.495	0.000	15091	0.25000	0.240
7 CROTONAL	6.408	6.408	0.000	13736	0.25000	0.241
9 BUTYRALS/MEK	7.061	7.061	0.000	12498	0.25000	0.233
10 BENZALDEHYDE	8.061	8.061	0.000	9441	0.25000	0.247
11 ISO-PENTANAL	9.148	9.148	0.000	10500	0.25000	0.232
12 PENTANAL	9.821	9.821	0.000	10423	0.25000	0.235
14 O-TOLUALDEHYDE	10.874	10.874	0.000	9638	0.25000	0.260
15 M,P-TOLUALDEHYDE	11.188	11.188	0.000	16415	0.50000	0.524
16 HEXANAL	12.768	12.768	0.000	11060	0.25000	0.269
17 2,5-Dimethylbenzaldehyde	13.354	13.354	0.000	8339	0.25000	0.274

Data File: /chem/hp1cf.i/14Sep2010.b/f0914016.d

Date : 14-SEP-2010 19:12

Client ID: 3

Sample Info: f1838-34-0.25:3

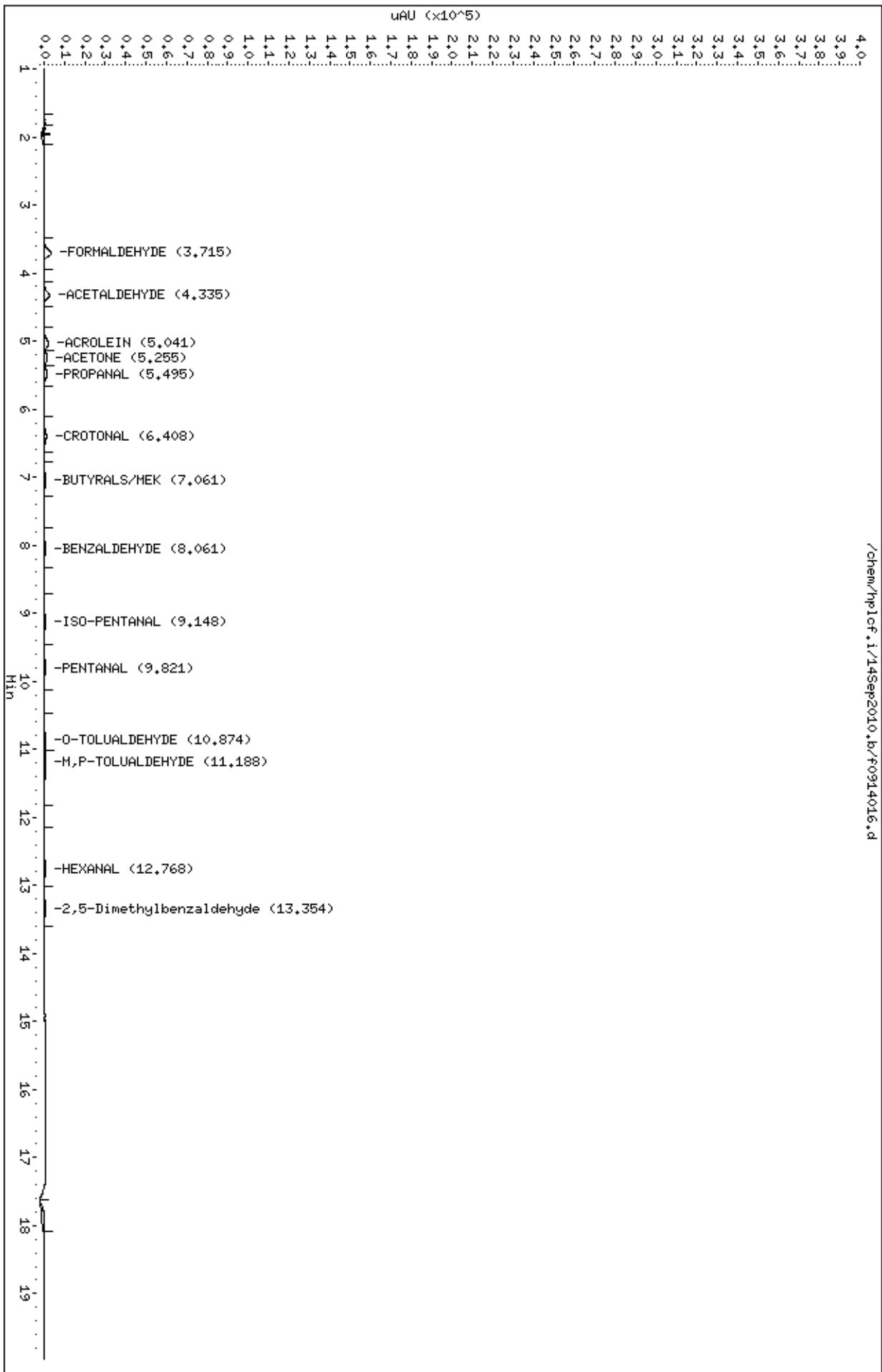
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
 Data file : /chem/hplcf.i/14Sep2010.b/f0914015.d
 Lab Smp Id: 1838-34-0.25 Client Smp ID: 3
 Inj Date : 14-SEP-2010 18:51
 Operator : CRL Inst ID: hplcf.i
 Smp Info : ;1838-34-0.25;3
 Misc Info :
 Comment :
 Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
 Meth Date : 15-Sep-2010 09:39 cleaf Quant Type: ESTD
 Cal Date : 14-SEP-2010 18:51 Cal File: f0914015.d
 Als bottle: 1 Calibration Sample, Level: 3
 Dil Factor: 1.00000
 Integrator: Falcon Compound Sublist: all.sub
 Target Version: 3.50
 Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
 Vs 5.00000 Sample Final Volume (mL)
 Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH	Compound Not Detected.					
2 FORMALDEHYDE	3.705	3.705	0.000	25860	0.25000	0.229
3 ACETALDEHYDE	4.325	4.325	0.000	20478	0.25000	0.240
4 ACROLEIN	5.025	5.025	0.000	19230	0.25000	0.234
5 ACETONE	5.238	5.238	0.000	15266	0.25000	0.242
6 PROPANAL	5.472	5.472	0.000	14945	0.25000	0.236
7 CROTONAL	6.385	6.385	0.000	13680	0.25000	0.238
9 BUTYRALS/MEK	7.032	7.032	0.000	12782	0.25000	0.234
10 BENZALDEHYDE	8.005	8.005	0.000	8867	0.25000	0.231
11 ISO-PENTANAL	9.085	9.085	0.000	11141	0.25000	0.241
12 PENTANAL	9.745	9.745	0.000	10552	0.25000	0.235
14 O-TOLUALDEHYDE	10.865	10.865	0.000	9186	0.25000	0.250
15 M,P-TOLUALDEHYDE	11.165	11.165	0.000	15167	0.50000	0.490
16 HEXANAL	12.725	12.725	0.000	9658	0.25000	0.240
17 2,5-Dimethylbenzaldehyde	13.318	13.318	0.000	7658	0.25000	0.258

Data File: /chem/hp1cf.i/14Sep2010.b/f0914015.d

Date : 14-SEP-2010 18:51

Client ID: 3

Sample Info: f1838-34-0.25;3

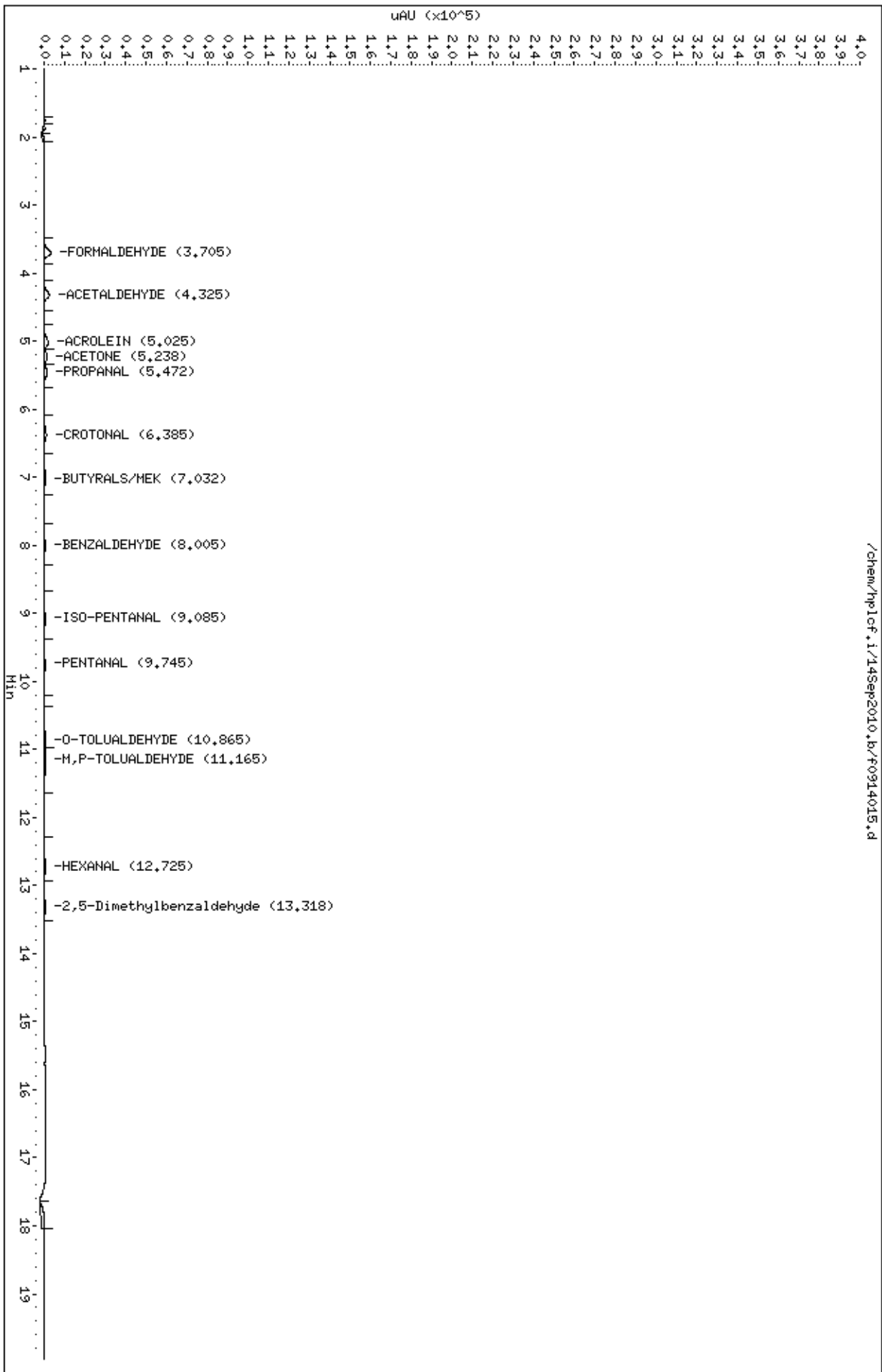
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914020.d
Lab Smp Id: 1838-34-0.5 Client Smp ID: 4
Inj Date : 14-SEP-2010 20:35
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-0.5;4
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:39 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 20:35 Cal File: f0914020.d
Als bottle: 1 Calibration Sample, Level: 4
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH	Compound Not Detected.					
2 FORMALDEHYDE	3.745	3.745	0.000	44981	0.50000	0.420
3 ACETALDEHYDE	4.378	4.378	0.000	42208	0.50000	0.499
4 ACRROLEIN	5.098	5.098	0.000	40725	0.50000	0.505
5 ACETONE	5.305	5.305	0.000	31112	0.50000	0.498
6 PROPANAL	5.545	5.545	0.000	31593	0.50000	0.505
7 CROTONAL	6.492	6.492	0.000	28714	0.50000	0.501
9 BUTYRALS/MEK	7.159	7.159	0.000	26645	0.50000	0.495
10 BENZALDEHYDE	8.165	8.165	0.000	18778	0.50000	0.491
11 ISO-PENTANAL	9.285	9.285	0.000	22133	0.50000	0.490
12 PENTANAL	9.979	9.979	0.000	22290	0.50000	0.501
14 O-TOLUALDEHYDE	11.039	11.039	0.000	18580	0.50000	0.501
15 M,P-TOLUALDEHYDE	11.338	11.338	0.000	30599	1.00000	0.980
16 HEXANAL	12.872	12.872	0.000	20818	0.50000	0.503
17 2,5-Dimethylbenzaldehyde	13.465	13.465	0.000	15377	0.50000	0.494

Data File: /chem/hp1cf.i/14Sep2010.b/f0914020.d

Date : 14-SEP-2010 20:35

Client ID: 4

Sample Info: f1838-34-0.5;4

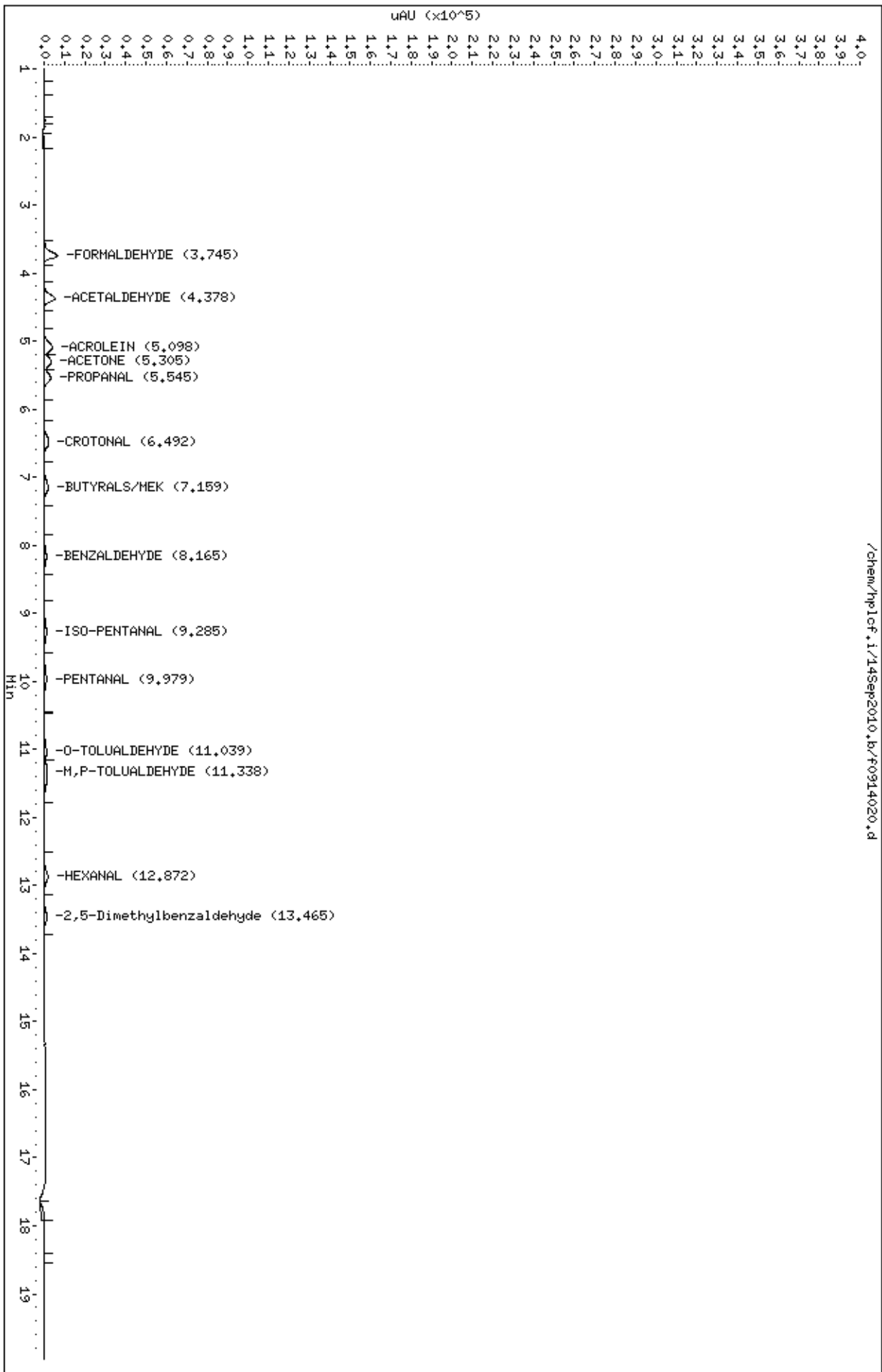
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914019.d
Lab Smp Id: 1838-34-0.5 Client Smp ID: 4
Inj Date : 14-SEP-2010 20:15
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-0.5;4
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:39 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 20:15 Cal File: f0914019.d
Als bottle: 1 Calibration Sample, Level: 4
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH	Compound Not Detected.					
2 FORMALDEHYDE	3.734	3.734	0.000	45340	0.50000	0.419
3 ACETALDEHYDE	4.367	4.367	0.000	41778	0.50000	0.494
4 ACRROLEIN	5.081	5.081	0.000	40776	0.50000	0.507
5 ACETONE	5.287	5.287	0.000	31381	0.50000	0.503
6 PROPANAL	5.534	5.534	0.000	31039	0.50000	0.497
7 CROTONAL	6.468	6.468	0.000	29363	0.50000	0.513
9 BUTYRALS/MEK	7.128	7.128	0.000	28301	0.50000	0.525
10 BENZALDEHYDE	8.141	8.141	0.000	18877	0.50000	0.493
11 ISO-PENTANAL	9.241	9.241	0.000	22567	0.50000	0.498
12 PENTANAL	9.921	9.921	0.000	22543	0.50000	0.507
14 O-TOLUALDEHYDE	10.988	10.988	0.000	18387	0.50000	0.496
15 M,P-TOLUALDEHYDE	11.307	11.307	0.000	31335	1.00000	1.00
16 HEXANAL	12.848	12.848	0.000	20238	0.50000	0.489
17 2,5-Dimethylbenzaldehyde	13.434	13.434	0.000	15337	0.50000	0.493

Data File: /chem/hp1cf.i/14Sep2010.b/f0914019.d

Date : 14-SEP-2010 20:15

Client ID: 4

Sample Info: f1838-34-0.5;4

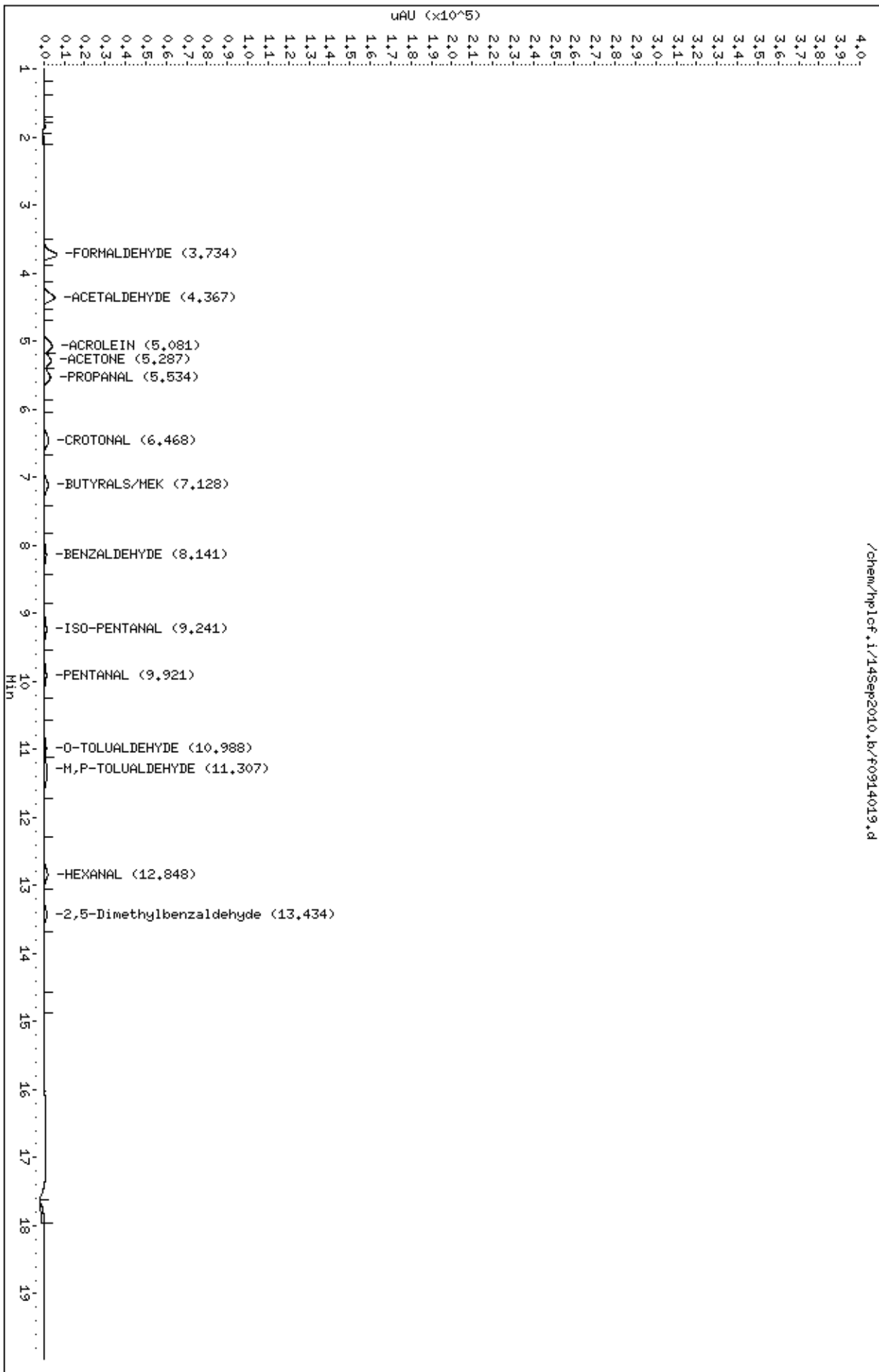
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914018.d
Lab Smp Id: 1838-34-0.5 Client Smp ID: 4
Inj Date : 14-SEP-2010 19:54
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-0.5;4
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:39 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 19:54 Cal File: f0914018.d
Als bottle: 1 Calibration Sample, Level: 4
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH	Compound Not Detected.					
2 FORMALDEHYDE	3.734	3.734	0.000	45518	0.50000	0.416
3 ACETALDEHYDE	4.361	4.361	0.000	42551	0.50000	0.502
4 ACRROLEIN	5.081	5.081	0.000	40341	0.50000	0.502
5 ACETONE	5.287	5.287	0.000	31502	0.50000	0.505
6 PROPANAL	5.534	5.534	0.000	30923	0.50000	0.495
7 CROTONAL	6.467	6.467	0.000	28980	0.50000	0.508
9 BUTYRALS/MEK	7.127	7.127	0.000	27428	0.50000	0.512
10 BENZALDEHYDE	8.127	8.127	0.000	19085	0.50000	0.497
11 ISO-PENTANAL	9.227	9.227	0.000	22964	0.50000	0.506
12 PENTANAL	9.907	9.907	0.000	22746	0.50000	0.512
14 O-TOLUALDEHYDE	10.987	10.987	0.000	18857	0.50000	0.508
15 M,P-TOLUALDEHYDE	11.307	11.307	0.000	31298	1.00000	1.000
16 HEXANAL	12.834	12.834	0.000	20952	0.50000	0.505
17 2,5-Dimethylbenzaldehyde	13.421	13.421	0.000	15817	0.50000	0.507

Data File: /chem/hp1cf.i/14Sep2010.b/f0914018.d

Date : 14-SEP-2010 19:54

Client ID: 4

Sample Info: f1838-34-0.5;4

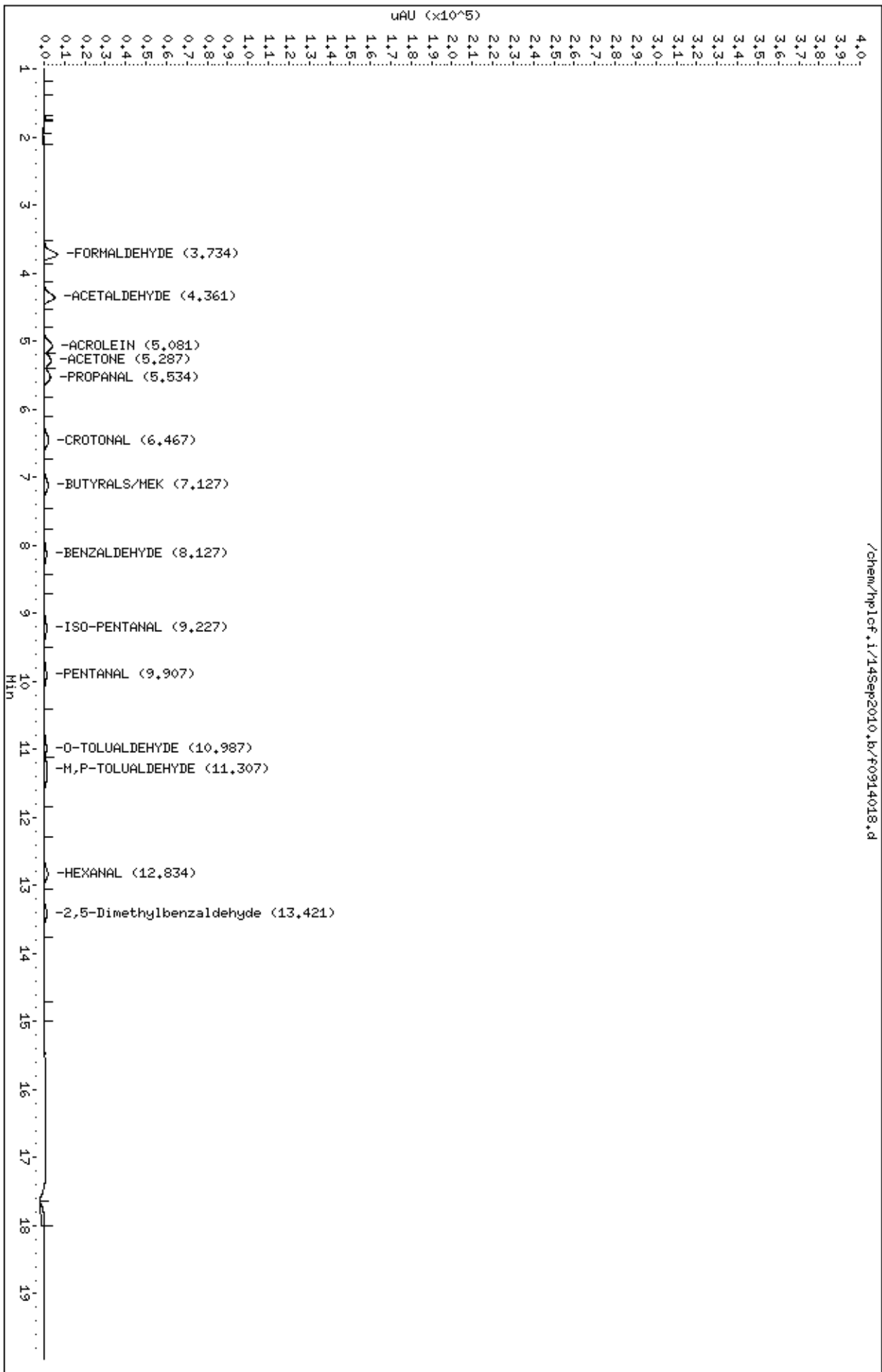
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914023.d
Lab Smp Id: 1838-34-1.0 Client Smp ID: 5
Inj Date : 14-SEP-2010 21:38
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-1.0;5
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 21:38 Cal File: f0914023.d
Als bottle: 1 Calibration Sample, Level: 5
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH	Compound Not Detected.					
2 FORMALDEHYDE	3.739	3.739	0.000	105489	1.00000	0.988
3 ACETALDEHYDE	4.365	4.365	0.000	82932	1.00000	0.984
4 ACRROLEIN	5.085	5.085	0.000	78338	1.00000	0.980
5 ACETONE	5.292	5.292	0.000	60890	1.00000	0.981
6 PROPANAL	5.539	5.539	0.000	61135	1.00000	0.983
7 CROTONAL	6.472	6.472	0.000	55126	1.00000	0.971
9 BUTYRALS/MEK	7.139	7.139	0.000	52870	1.00000	0.986
10 BENZALDEHYDE	8.159	8.159	0.000	36794	1.00000	0.970
11 ISO-PENTANAL	9.265	9.265	0.000	44316	1.00000	0.987
12 PENTANAL	9.939	9.939	0.000	43634	1.00000	0.987
14 O-TOLUALDEHYDE	11.019	11.019	0.000	35016	1.00000	0.957
15 M,P-TOLUALDEHYDE	11.325	11.325	0.000	59329	2.00000	1.93
16 HEXANAL	12.852	12.852	0.000	39118	1.00000	0.957
17 2,5-Dimethylbenzaldehyde	13.445	13.445	0.000	28339	1.00000	0.929

Data File: /chem/hp1cf.i/14Sep2010.b/f0914023.d

Date : 14-SEP-2010 21:38

Client ID: 5

Sample Info: f1838-34-1.0;5

Purge Volume: 5.0

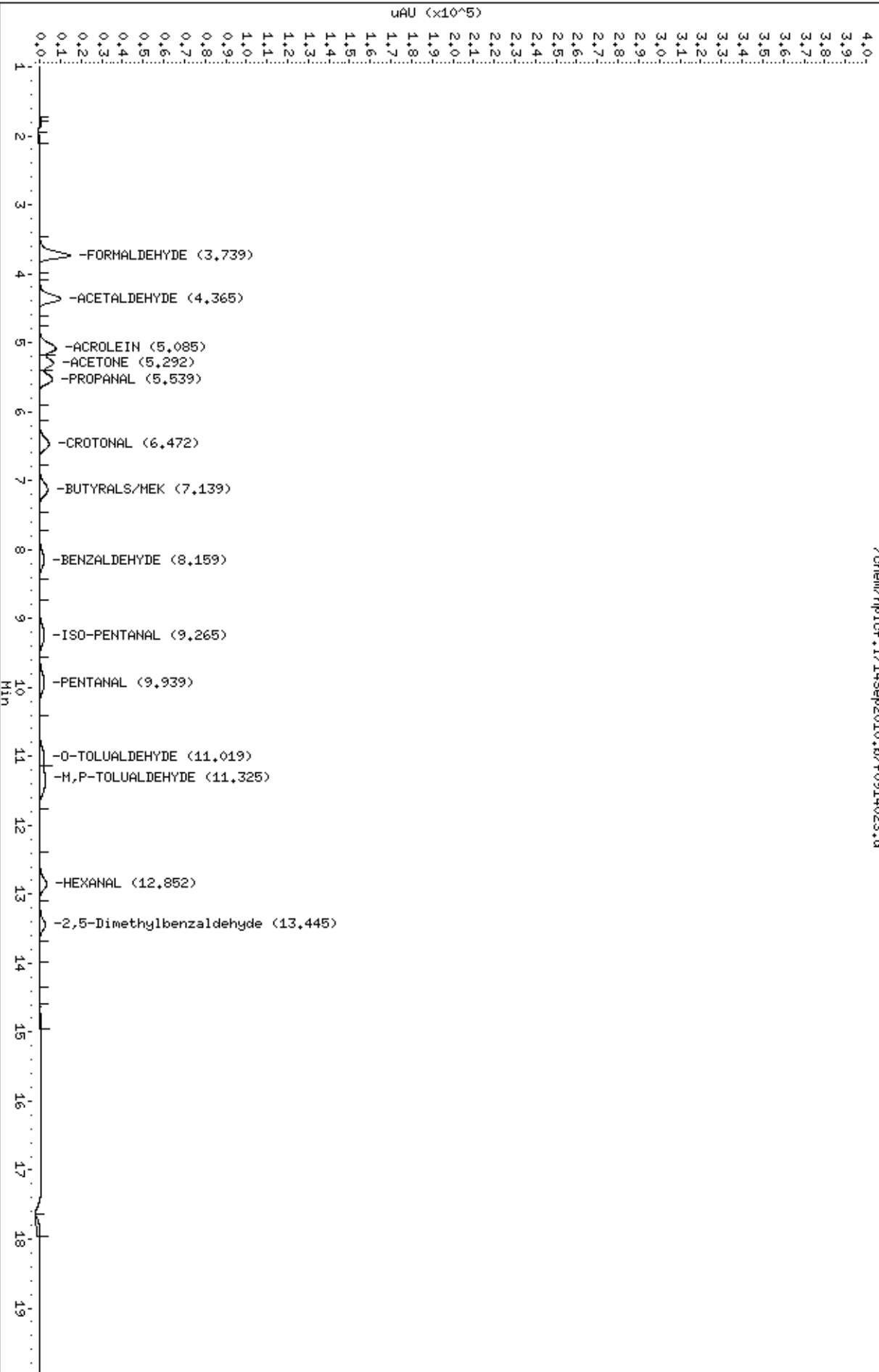
Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00

/chem/hp1cf.i/14Sep2010.b/f0914023.d



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
 Data file : /chem/hplcf.i/14Sep2010.b/f0914022.d
 Lab Smp Id: 1838-34-1.0 Client Smp ID: 5
 Inj Date : 14-SEP-2010 21:17
 Operator : CRL Inst ID: hplcf.i
 Smp Info : ;1838-34-1.0;5
 Misc Info :
 Comment :
 Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
 Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
 Cal Date : 14-SEP-2010 21:17 Cal File: f0914022.d
 Als bottle: 1 Calibration Sample, Level: 5
 Dil Factor: 1.00000
 Integrator: Falcon Compound Sublist: all.sub
 Target Version: 3.50
 Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
 Vs 5.00000 Sample Final Volume (mL)
 Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH	Compound Not Detected.					
2 FORMALDEHYDE	3.747	3.747	0.000	105394	1.00000	0.987
3 ACETALDEHYDE	4.373	4.373	0.000	83089	1.00000	0.985
4 ACRROLEIN	5.093	5.093	0.000	77739	1.00000	0.971
5 ACETONE	5.306	5.306	0.000	61069	1.00000	0.982
6 PROPANAL	5.546	5.546	0.000	61304	1.00000	0.985
7 CROTONAL	6.487	6.487	0.000	55721	1.00000	0.979
9 BUTYRALS/MEK	7.153	7.153	0.000	52676	1.00000	0.981
10 BENZALDEHYDE	8.180	8.180	0.000	37230	1.00000	0.979
11 ISO-PENTANAL	9.273	9.273	0.000	43396	1.00000	0.965
12 PENTANAL	9.960	9.960	0.000	42834	1.00000	0.968
14 O-TOLUALDEHYDE	11.027	11.027	0.000	35079	1.00000	0.955
15 M,P-TOLUALDEHYDE	11.340	11.340	0.000	58406	2.00000	1.89
16 HEXANAL	12.867	12.867	0.000	39102	1.00000	0.952
17 2,5-Dimethylbenzaldehyde	13.453	13.453	0.000	29132	1.00000	0.948

Data File: /chem/hp1cf.i/14Sep2010.b/f0914022.d

Date : 14-SEP-2010 21:17

Client ID: 5

Sample Info: #1838-34-1.0;5

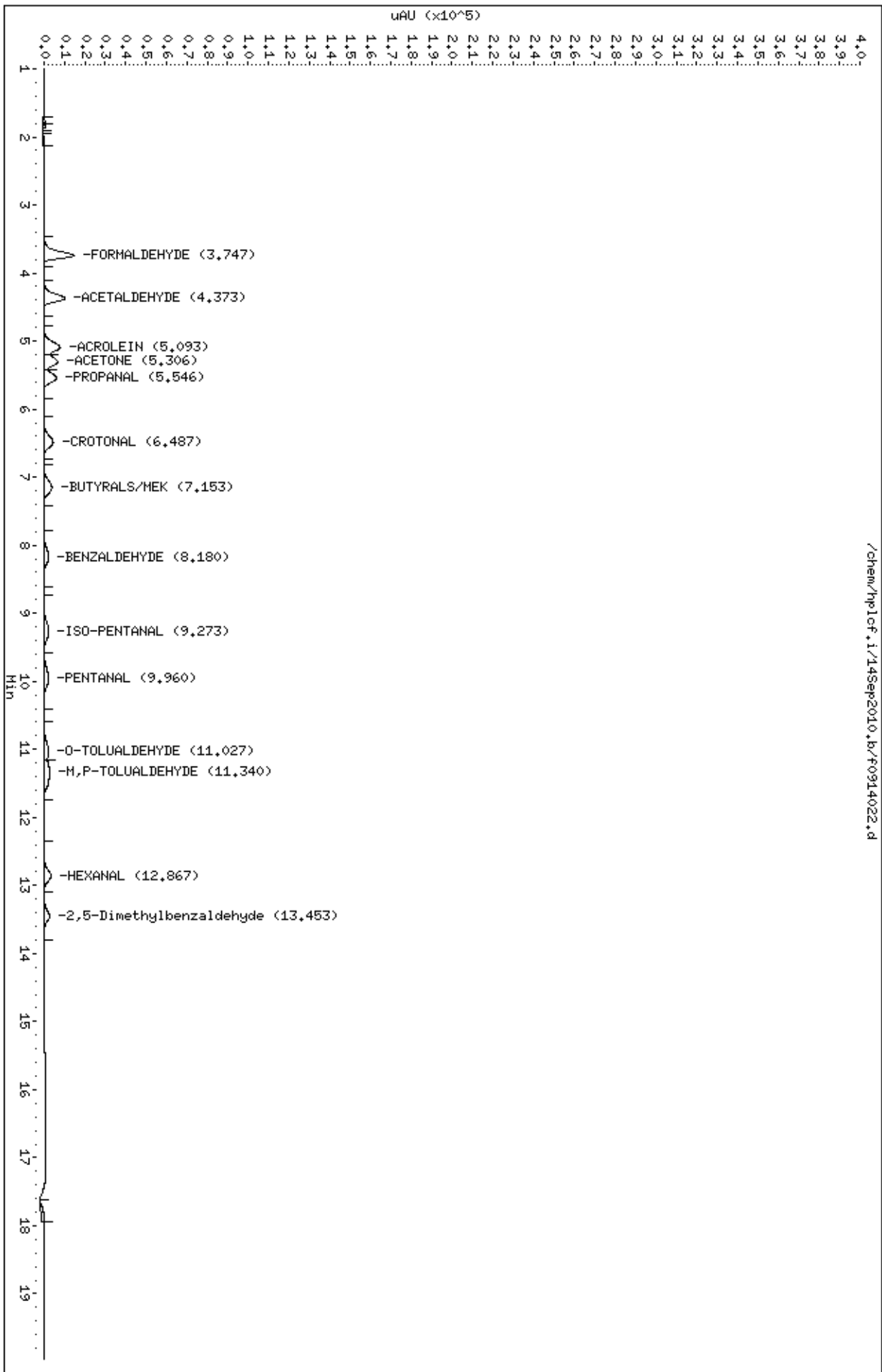
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914021.d
Lab Smp Id: 1838-34-1.0 Client Smp ID: 5
Inj Date : 14-SEP-2010 20:56
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-1.0;5
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 20:56 Cal File: f0914021.d
Als bottle: 1 Calibration Sample, Level: 5
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
1 DNPH	Compound Not Detected.					
2 FORMALDEHYDE	3.740	3.740	0.000	105805	1.00000	0.990
3 ACETALDEHYDE	4.374	4.374	0.000	82817	1.00000	0.981
4 ACRROLEIN	5.094	5.094	0.000	77980	1.00000	0.971
5 ACETONE	5.301	5.301	0.000	60932	1.00000	0.979
6 PROPANAL	5.547	5.547	0.000	60789	1.00000	0.975
7 CROTONAL	6.487	6.487	0.000	55018	1.00000	0.964
9 BUTYRALS/MEK	7.147	7.147	0.000	53240	1.00000	0.990
10 BENZALDEHYDE	8.161	8.161	0.000	36832	1.00000	0.967
11 ISO-PENTANAL	9.274	9.274	0.000	44231	1.00000	0.980
12 PENTANAL	9.960	9.960	0.000	43614	1.00000	0.982
14 O-TOLUALDEHYDE	11.021	11.021	0.000	35206	1.00000	0.954
15 M,P-TOLUALDEHYDE	11.334	11.334	0.000	59111	2.00000	1.90
16 HEXANAL	12.860	12.860	0.000	39766	1.00000	0.964
17 2,5-Dimethylbenzaldehyde	13.447	13.447	0.000	28909	1.00000	0.936

Data File: /chem/hp1cf.i/14Sep2010.b/f0914021.d

Date : 14-SEP-2010 20:56

Client ID: 5

Sample Info: #1838-34-1.0;5

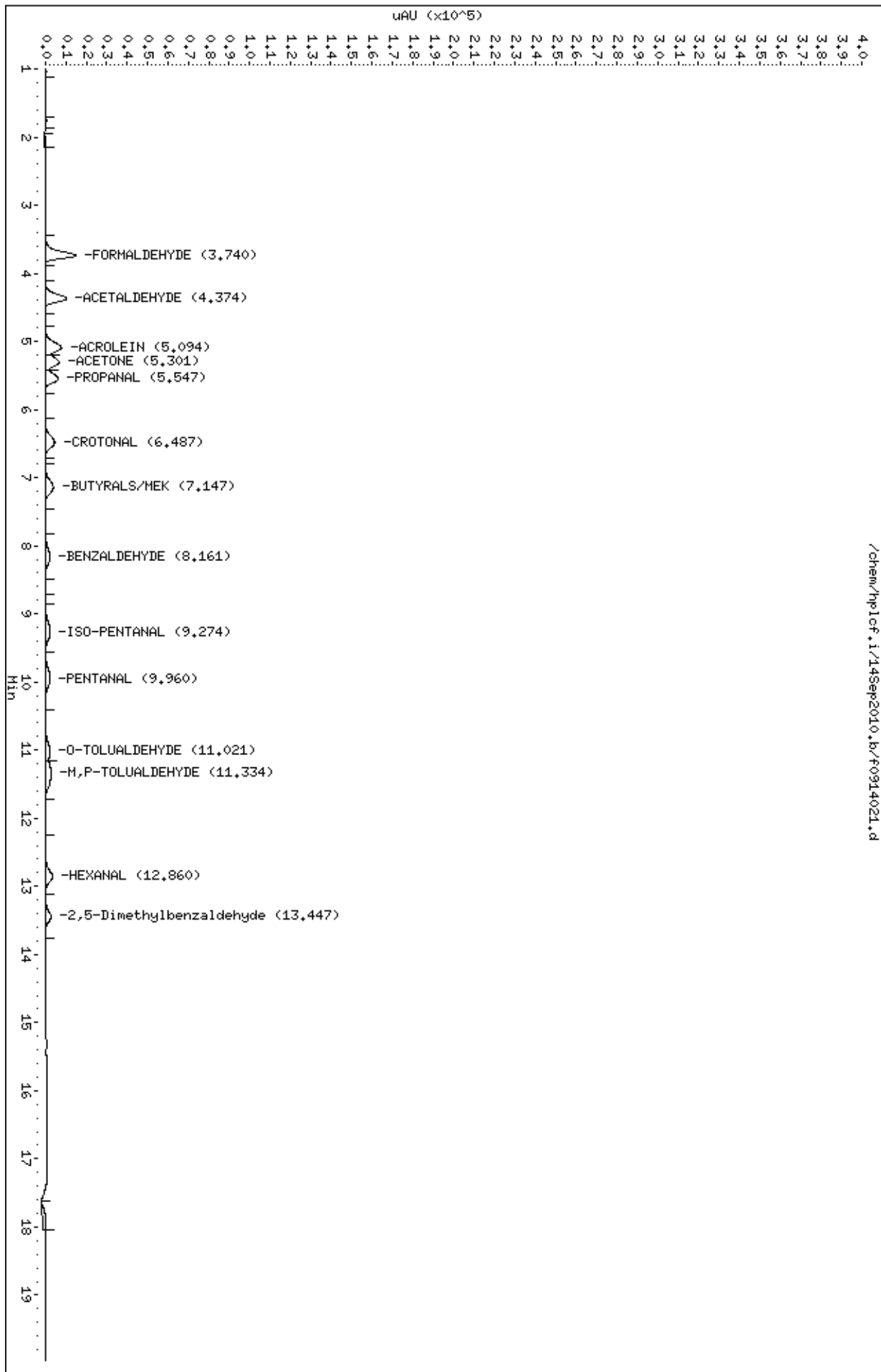
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914026.d
Lab Smp Id: 1838-32-4.0 Client Smp ID: 6
Inj Date : 14-SEP-2010 22:41
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-32-4.0;6
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:29 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 22:41 Cal File: f0914026.d
Als bottle: 1 Calibration Sample, Level: 6
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.737	3.737	0.000	450621	4.00000	4.01
3 ACETALDEHYDE	4.371	4.371	0.000	352416	4.00000	4.00
4 ACRROLEIN	5.091	5.091	0.000	335424	4.00000	4.00
5 ACETONE	5.298	5.298	0.000	258255	4.00000	4.04
6 PROPANAL	5.537	5.537	0.000	256942	4.00000	3.98
7 CROTONAL	6.478	6.478	0.000	233257	4.00000	4.00
9 BUTYRALS/MEK	7.144	7.144	0.000	222115	4.00000	3.99
10 BENZALDEHYDE	8.158	8.158	0.000	157668	4.00000	4.01
11 ISO-PENTANAL	9.258	9.258	0.000	187027	4.00000	4.00
12 PENTANAL	9.938	9.938	0.000	182923	4.00000	3.99
14 O-TOLUALDEHYDE	11.011	11.011	0.000	147236	4.00000	4.02
15 M,P-TOLUALDEHYDE	11.324	11.324	0.000	248923	8.00000	7.99
16 HEXANAL	12.851	12.851	0.000	163558	4.00000	4.01
17 2,5-Dimethylbenzaldehyde	13.444	13.444	0.000	117296	4.00000	3.99

Data File: /chem/hp1cf.i/14Sep2010.b/f0914026.d

Date : 14-SEP-2010 22:41

Client ID: 6

Sample Info: f1838-32-4.0;6

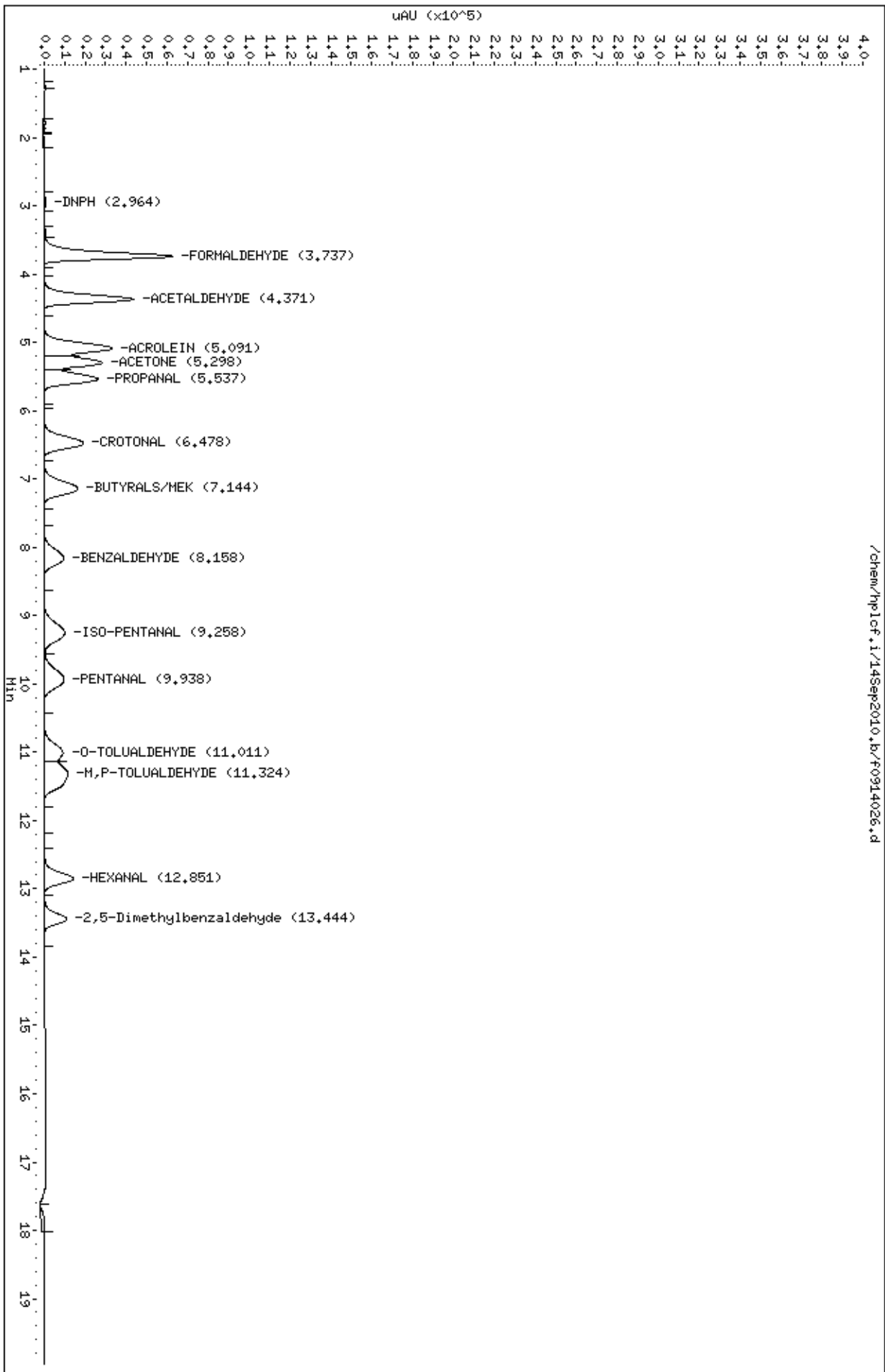
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914025.d
Lab Smp Id: 1838-32-4.0 Client Smp ID: 6
Inj Date : 14-SEP-2010 22:20
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-32-4.0;6
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:29 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 22:20 Cal File: f0914025.d
Als bottle: 1 Calibration Sample, Level: 6
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.742	3.742	0.000	449099	4.00000	4.00
3 ACETALDEHYDE	4.375	4.375	0.000	352551	4.00000	4.01
4 ACRROLEIN	5.095	5.095	0.000	334904	4.00000	4.00
5 ACETONE	5.302	5.302	0.000	254054	4.00000	4.00
6 PROPANAL	5.542	5.542	0.000	258747	4.00000	3.99
7 CROTONAL	6.482	6.482	0.000	233603	4.00000	4.01
9 BUTYRALS/MEK	7.142	7.142	0.000	223717	4.00000	4.02
10 BENZALDEHYDE	8.162	8.162	0.000	157371	4.00000	4.01
11 ISO-PENTANAL	9.262	9.262	0.000	185991	4.00000	3.98
12 PENTANAL	9.942	9.942	0.000	182460	4.00000	3.97
14 O-TOLUALDEHYDE	11.009	11.009	0.000	145838	4.00000	3.99
15 M,P-TOLUALDEHYDE	11.322	11.322	0.000	249513	8.00000	8.01
16 HEXANAL	12.849	12.849	0.000	163297	4.00000	4.00
17 2,5-Dimethylbenzaldehyde	13.442	13.442	0.000	117509	4.00000	3.99

Data File: /chem/hp1cf.i/14Sep2010.b/f0914025.d

Date : 14-SEP-2010 22:20

Client ID: 6

Sample Info: f1838-32-4.0;6

Purge Volume: 5.0

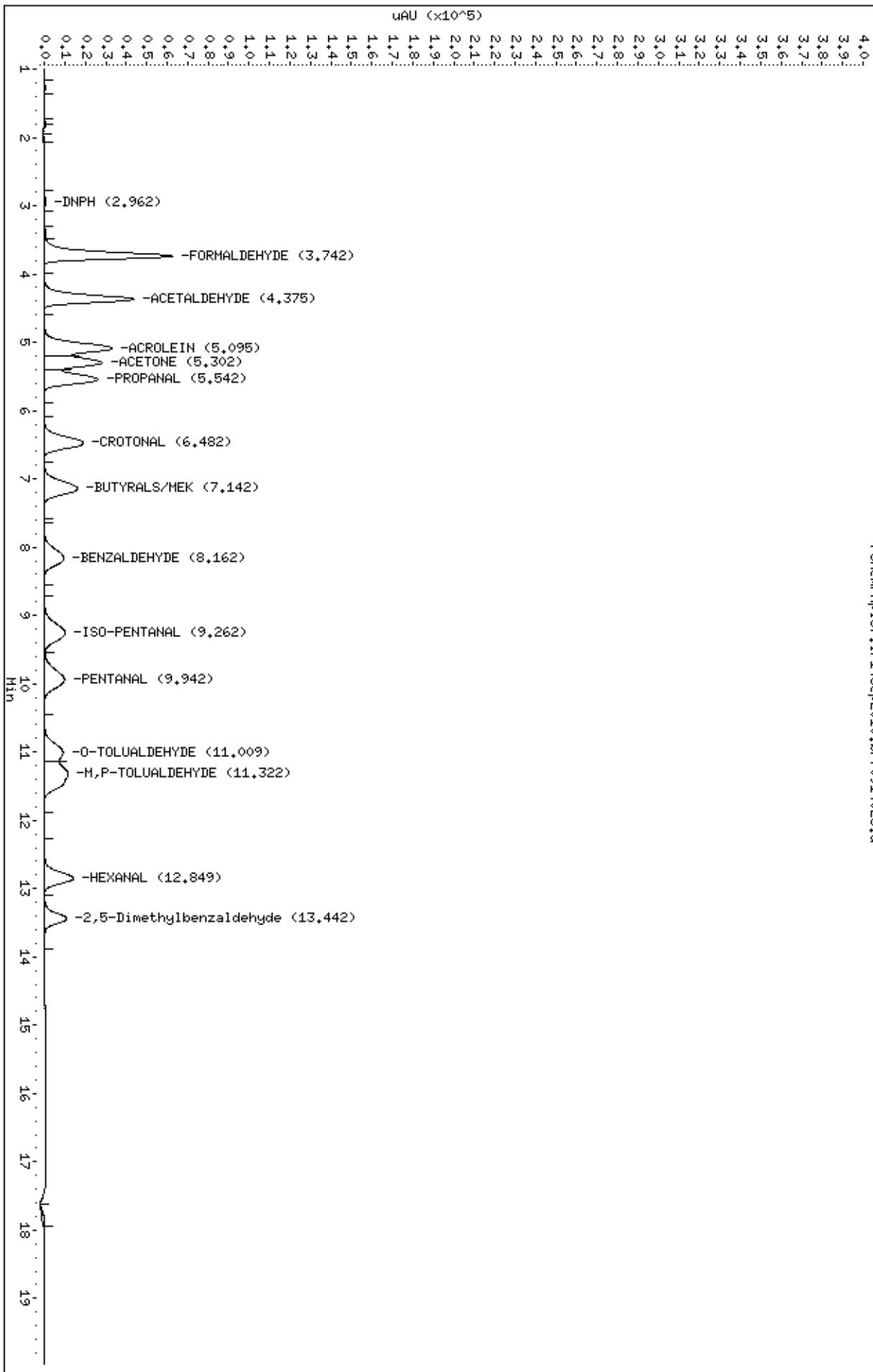
Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00

/chem/hp1cf.i/14Sep2010.b/f0914025.d



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
 Data file : /chem/hplcf.i/14Sep2010.b/f0914024.d
 Lab Smp Id: 1838-32-4.0 Client Smp ID: 6
 Inj Date : 14-SEP-2010 21:59
 Operator : CRL Inst ID: hplcf.i
 Smp Info : ;1838-32-4.0;6
 Misc Info :
 Comment :
 Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
 Meth Date : 15-Sep-2010 09:29 cleaf Quant Type: ESTD
 Cal Date : 14-SEP-2010 21:59 Cal File: f0914024.d
 Als bottle: 1 Calibration Sample, Level: 6
 Dil Factor: 1.00000
 Integrator: Falcon Compound Sublist: all.sub
 Target Version: 3.50
 Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
 Vs 5.00000 Sample Final Volume (mL)
 Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.742	3.742	0.000	448858	4.00000	4.00
3 ACETALDEHYDE	4.375	4.375	0.000	351227	4.00000	4.00
4 ACRROLEIN	5.096	5.096	0.000	334499	4.00000	4.00
5 ACETONE	5.302	5.302	0.000	254555	4.00000	4.00
6 PROPANAL	5.542	5.542	0.000	259688	4.00000	4.00
7 CROTONAL	6.482	6.482	0.000	232736	4.00000	4.00
9 BUTYRALS/MEK	7.142	7.142	0.000	221997	4.00000	4.00
10 BENZALDEHYDE	8.162	8.162	0.000	156893	4.00000	4.00
11 ISO-PENTANAL	9.269	9.269	0.000	187764	4.00000	4.00
12 PENTANAL	9.956	9.956	0.000	184847	4.00000	4.00
14 O-TOLUALDEHYDE	11.022	11.022	0.000	146742	4.00000	4.00
15 M,P-TOLUALDEHYDE	11.335	11.335	0.000	249029	8.00000	8.00
16 HEXANAL	12.855	12.855	0.000	163040	4.00000	4.00
17 2,5-Dimethylbenzaldehyde	13.449	13.449	0.000	118085	4.00000	4.00

Data File: /chem/hp1cf.i/14Sep2010.b/f0914024.d

Date : 14-SEP-2010 21:59

Client ID: 6

Sample Info: f1838-32-4.0;6

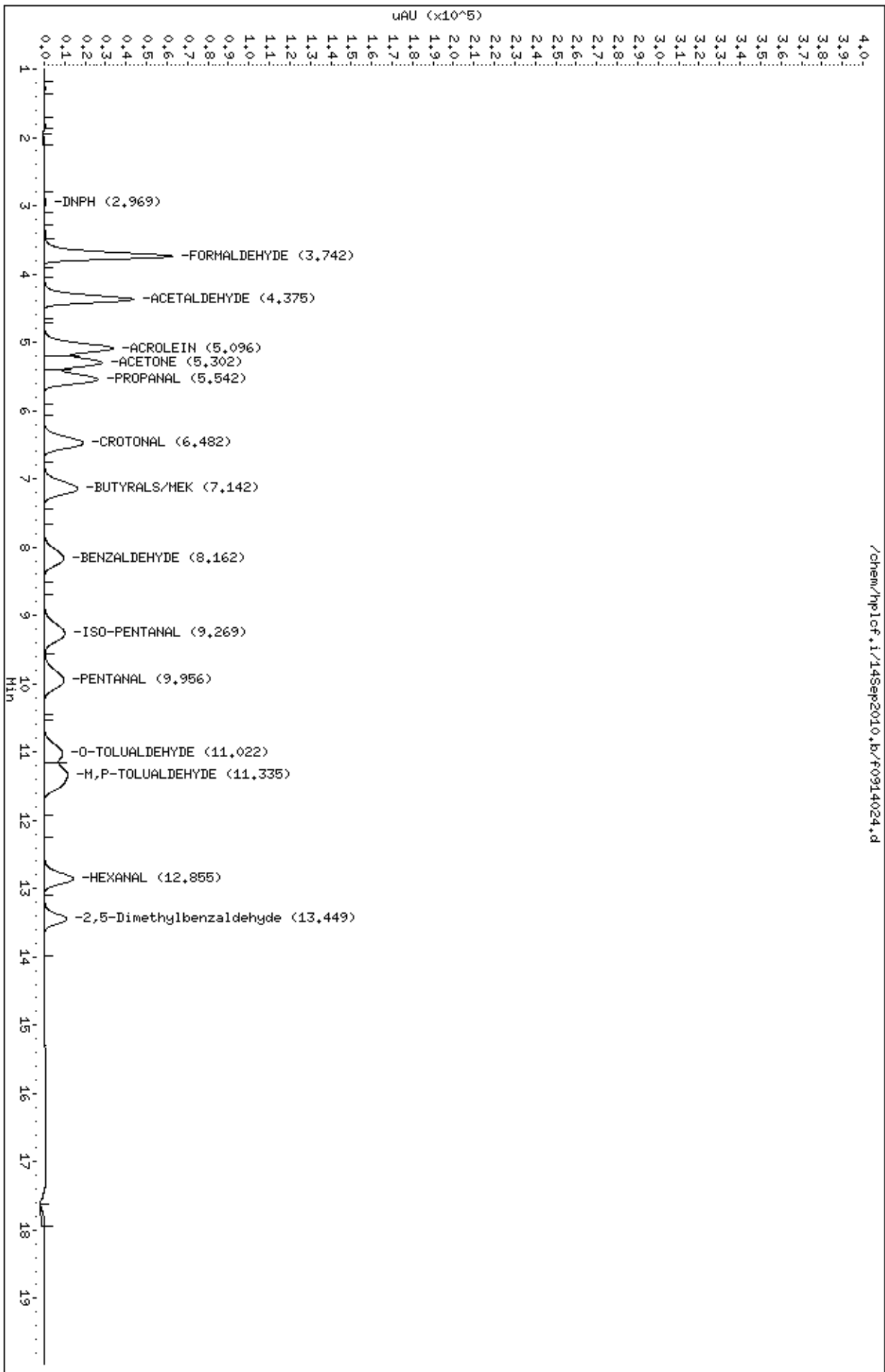
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914029.d
Lab Smp Id: 1838-34-10.0 Client Smp ID: 7
Inj Date : 14-SEP-2010 23:44
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-10.0;7
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 23:44 Cal File: f0914029.d
Als bottle: 1 Calibration Sample, Level: 7
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.737	3.737	0.000	1086878	10.0000	10.2
3 ACETALDEHYDE	4.370	4.370	0.000	830306	10.0000	9.88
4 ACRROLEIN	5.090	5.090	0.000	787034	10.0000	9.87
5 ACETONE	5.297	5.297	0.000	598880	10.0000	9.71
6 PROPANAL	5.537	5.537	0.000	611548	10.0000	9.88
7 CROTONAL	6.477	6.477	0.000	548148	10.0000	9.72
9 BUTYRALS/MEK	7.137	7.137	0.000	524094	10.0000	9.82
10 BENZALDEHYDE	8.157	8.157	0.000	371534	10.0000	9.84
11 ISO-PENTANAL	9.250	9.250	0.000	438929	10.0000	9.82
12 PENTANAL	9.937	9.937	0.000	430061	10.0000	9.78
14 O-TOLUALDEHYDE	11.010	11.010	0.000	347525	10.0000	9.59
15 M,P-TOLUALDEHYDE	11.323	11.323	0.000	585343	20.0000	19.2
16 HEXANAL	12.850	12.850	0.000	383168	10.0000	9.49
17 2,5-Dimethylbenzaldehyde	13.443	13.443	0.000	275222	10.0000	9.20

Data File: /chem/hp1cf.i/14Sep2010.b/f0914029.d

Date : 14-SEP-2010 23:44

Client ID: 7

Sample Info: #1838-34-10.0:7

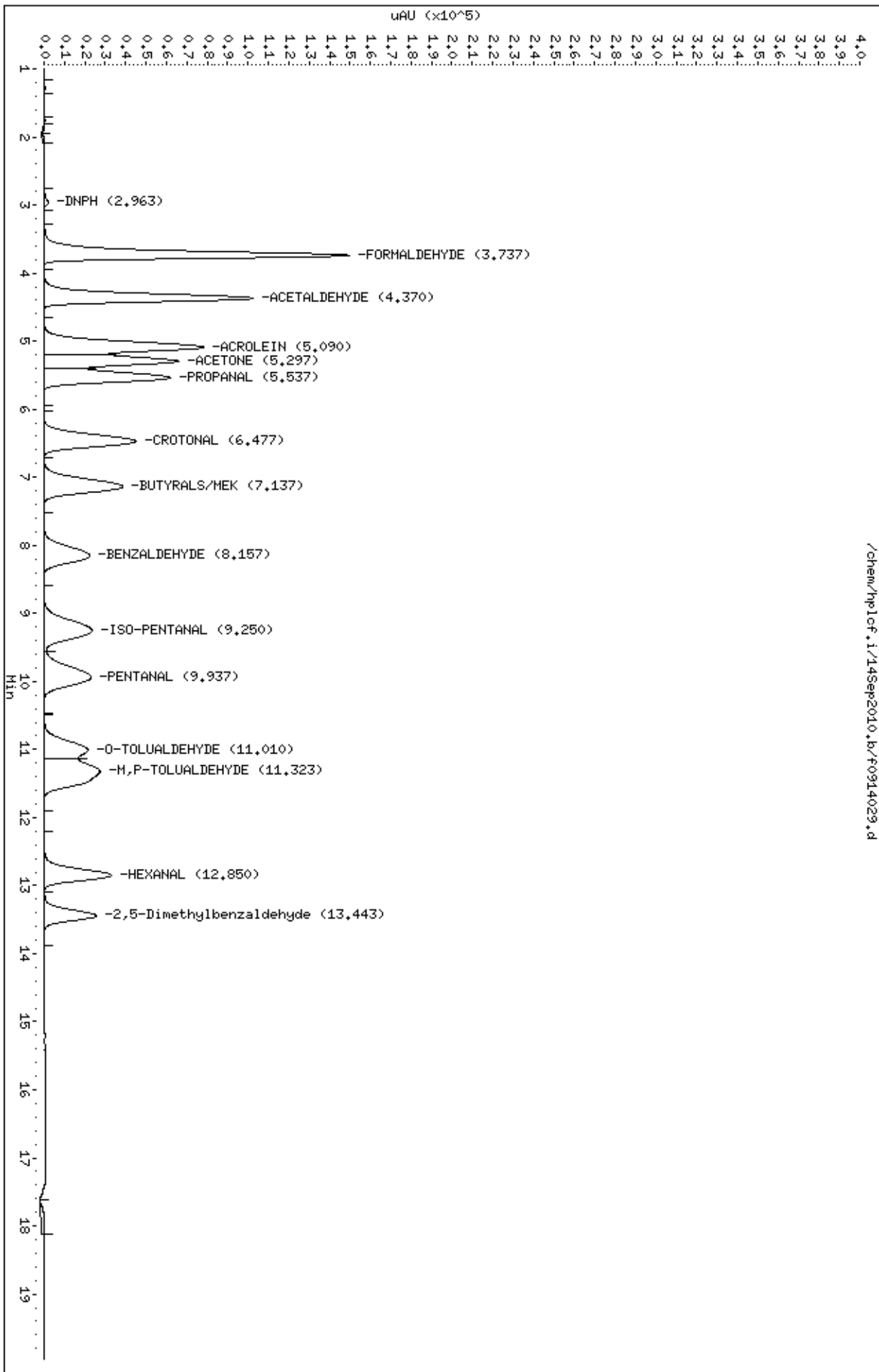
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914028.d
Lab Smp Id: 1838-34-10.0 Client Smp ID: 7
Inj Date : 14-SEP-2010 23:23
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-10.0;7
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 23:23 Cal File: f0914028.d
Als bottle: 1 Calibration Sample, Level: 7
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL - AMT (UG)	ON - COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.738	3.738	0.000	1087253	10.0000	10.2
3 ACETALDEHYDE	4.372	4.372	0.000	831049	10.0000	9.88
4 ACRROLEIN	5.092	5.092	0.000	792573	10.0000	9.93
5 ACETONE	5.298	5.298	0.000	597052	10.0000	9.66
6 PROPANAL	5.538	5.538	0.000	607745	10.0000	9.81
7 CROTONAL	6.478	6.478	0.000	548708	10.0000	9.71
9 BUTYRALS/MEK	7.138	7.138	0.000	524621	10.0000	9.81
10 BENZALDEHYDE	8.158	8.158	0.000	371558	10.0000	9.83
11 ISO-PENTANAL	9.258	9.258	0.000	439003	10.0000	9.81
12 PENTANAL	9.945	9.945	0.000	430505	10.0000	9.77
14 O-TOLUALDEHYDE	11.012	11.012	0.000	346558	10.0000	9.54
15 M,P-TOLUALDEHYDE	11.325	11.325	0.000	588014	20.0000	19.2
16 HEXANAL	12.852	12.852	0.000	383202	10.0000	9.46
17 2,5-Dimethylbenzaldehyde	13.445	13.445	0.000	275778	10.0000	9.16

Data File: /chem/hp1cf.i/14Sep2010.b/f0914028.d

Date : 14-SEP-2010 23:23

Client ID: 7

Sample Info: f1838-34-10.0:7

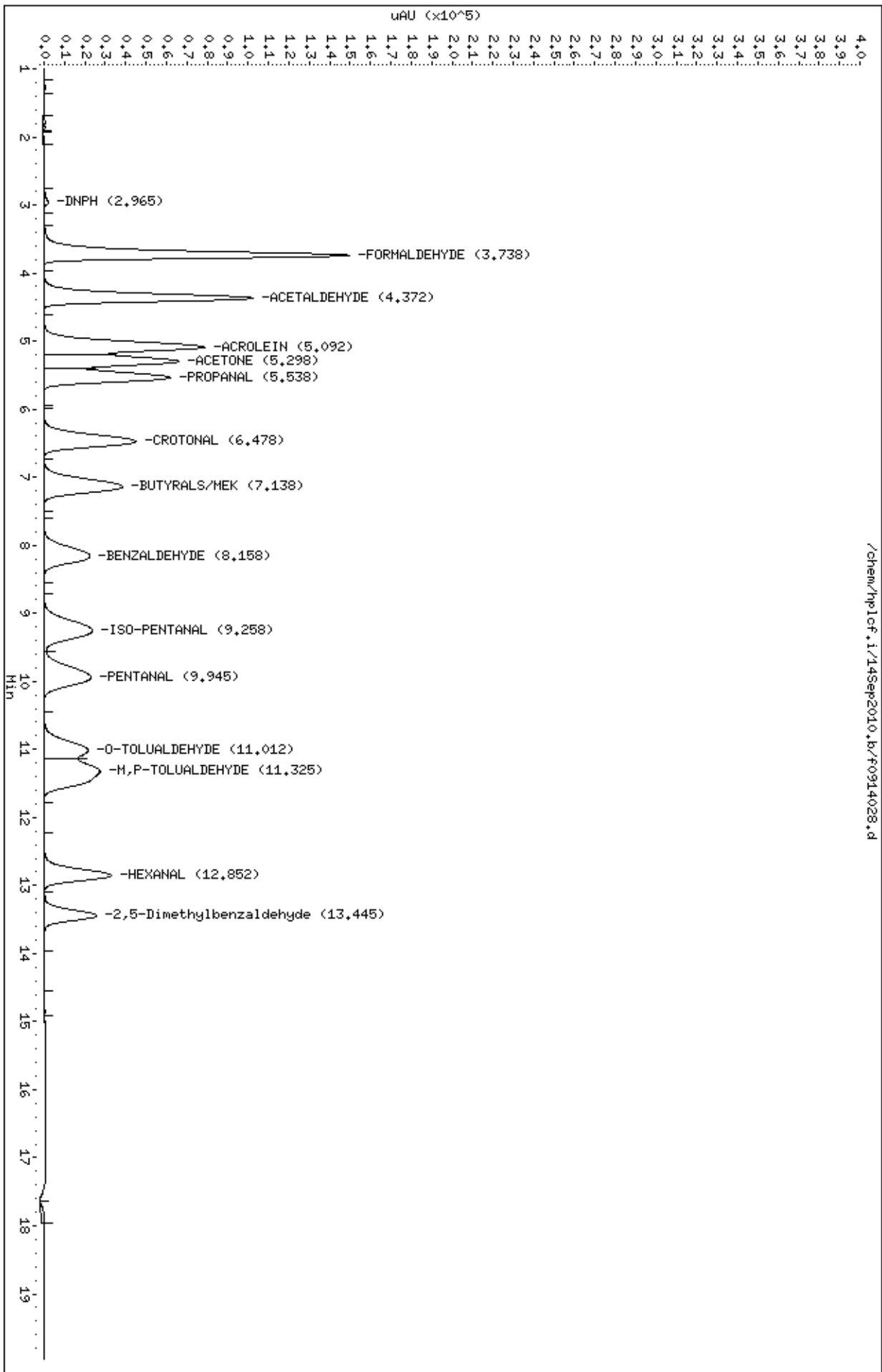
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914027.d
Lab Smp Id: 1838-34-10.0 Client Smp ID: 7
Inj Date : 14-SEP-2010 23:02
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-10.0;7
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 23:02 Cal File: f0914027.d
Als bottle: 1 Calibration Sample, Level: 7
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.738	3.738	0.000	1086932	10.0000	10.2
3 ACETALDEHYDE	4.371	4.371	0.000	830813	10.0000	9.87
4 ACRROLEIN	5.091	5.091	0.000	788842	10.0000	9.88
5 ACETONE	5.298	5.298	0.000	606307	10.0000	9.79
6 PROPANAL	5.544	5.544	0.000	602991	10.0000	9.72
7 CROTONAL	6.478	6.478	0.000	548541	10.0000	9.69
9 BUTYRALS/MEK	7.144	7.144	0.000	524696	10.0000	9.80
10 BENZALDEHYDE	8.164	8.164	0.000	371046	10.0000	9.80
11 ISO-PENTANAL	9.264	9.264	0.000	439896	10.0000	9.81
12 PENTANAL	9.944	9.944	0.000	430942	10.0000	9.76
14 O-TOLUALDEHYDE	11.018	11.018	0.000	348270	10.0000	9.55
15 M,P-TOLUALDEHYDE	11.331	11.331	0.000	585512	20.0000	19.1
16 HEXANAL	12.851	12.851	0.000	383719	10.0000	9.43
17 2,5-Dimethylbenzaldehyde	13.444	13.444	0.000	275329	10.0000	9.09

Data File: /chem/hp1cf.i/14Sep2010.b/f0914027.d

Date : 14-SEP-2010 23:02

Client ID: 7

Sample Info: #1838-34-10.0:7

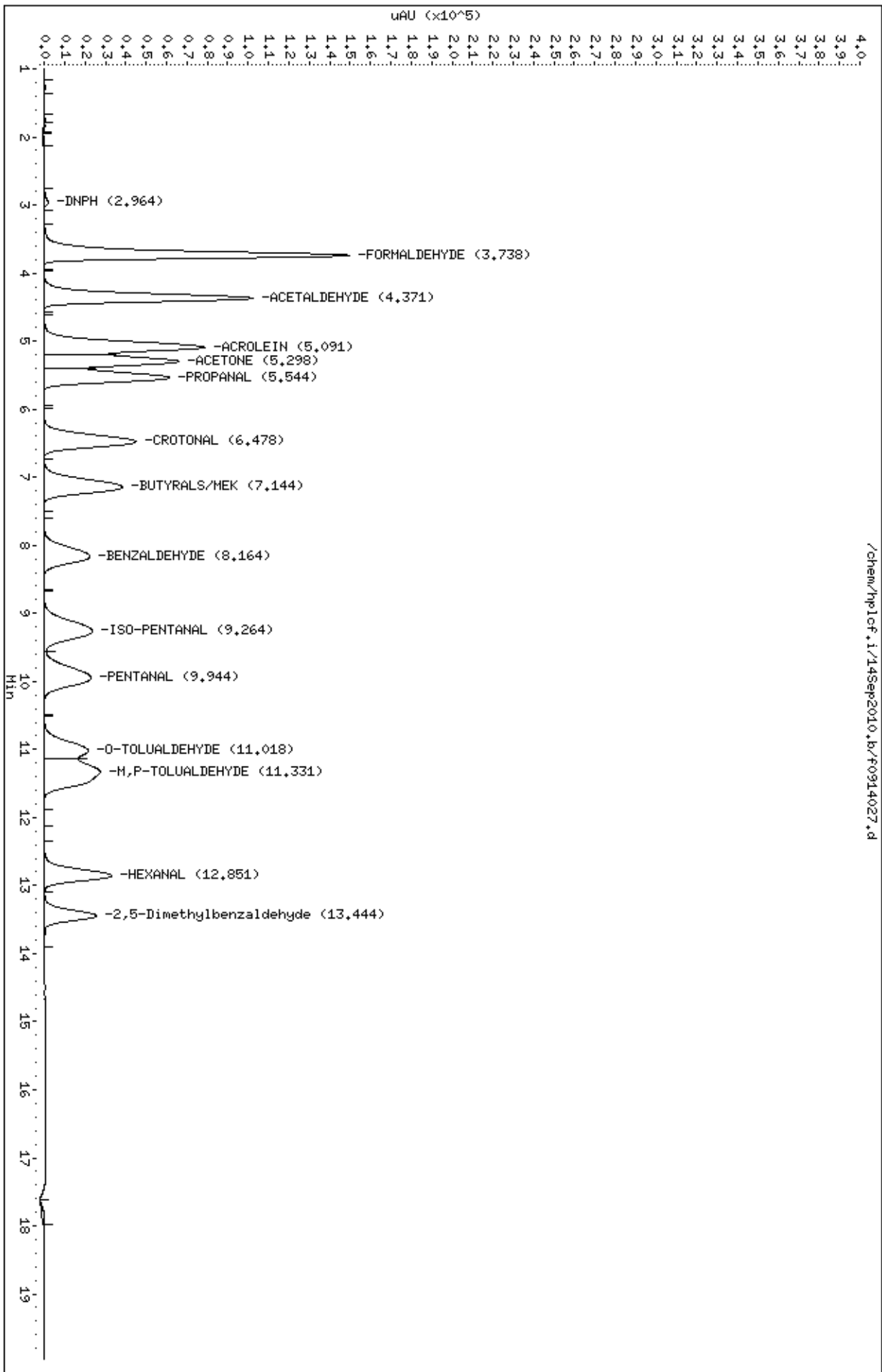
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914032.d
Lab Smp Id: 1838-34-20.0 Client Smp ID: 8
Inj Date : 15-SEP-2010 00:46
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-20.0;8
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 15-SEP-2010 00:46 Cal File: f0914032.d
Als bottle: 1 Calibration Sample, Level: 8
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.744	3.744	0.000	2254108	20.0000	20.9
3 ACETALDEHYDE	4.371	4.371	0.000	1715969	20.0000	20.4
4 ACRROLEIN	5.091	5.091	0.000	1648845	20.0000	20.6
5 ACETONE	5.297	5.297	0.000	1232915	20.0000	20.0
6 PROPANAL	5.544	5.544	0.000	1256858	20.0000	20.3
7 CROTONAL	6.484	6.484	0.000	1134806	20.0000	20.1
9 BUTYRALS/MEK	7.151	7.151	0.000	1085507	20.0000	20.3
10 BENZALDEHYDE	8.171	8.171	0.000	766541	20.0000	20.2
11 ISO-PENTANAL	9.271	9.271	0.000	915740	20.0000	20.4
12 PENTANAL	9.957	9.957	0.000	898851	20.0000	20.4
14 O-TOLUALDEHYDE	11.024	11.024	0.000	718581	20.0000	19.8
15 M,P-TOLUALDEHYDE	11.337	11.337	0.000	1214005	40.0000	39.9
16 HEXANAL	12.864	12.864	0.000	789706	20.0000	19.6
17 2,5-Dimethylbenzaldehyde	13.457	13.457	0.000	565128	20.0000	19.1

Data File: /chem/hplcf.i/14Sep2010.b/f0914032.d

Date : 15-SEP-2010 00:46

Client ID: 8

Sample Info: J1838-34-20.0:8

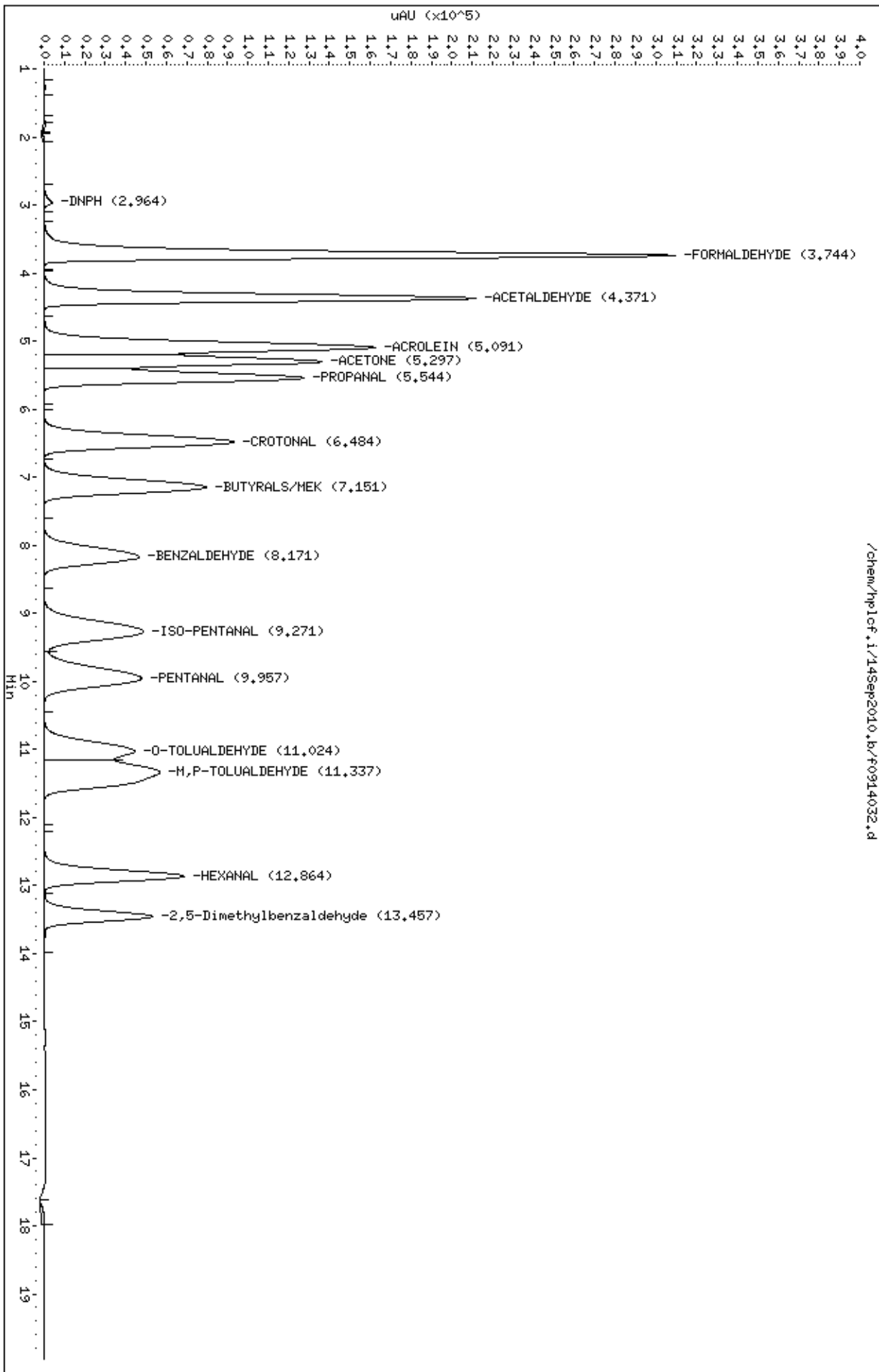
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914031.d
Lab Smp Id: 1838-34-20.0 Client Smp ID: 8
Inj Date : 15-SEP-2010 00:25
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-20.0;8
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 15-SEP-2010 00:25 Cal File: f0914031.d
Als bottle: 1 Calibration Sample, Level: 8
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.737	3.737	0.000	2253426	20.0000	21.0
3 ACETALDEHYDE	4.370	4.370	0.000	1715704	20.0000	20.4
4 ACRROLEIN	5.090	5.090	0.000	1623561	20.0000	20.3
5 ACETONE	5.297	5.297	0.000	1258750	20.0000	20.4
6 PROPANAL	5.544	5.544	0.000	1256634	20.0000	20.3
7 CROTONAL	6.484	6.484	0.000	1135155	20.0000	20.1
9 BUTYRALS/MEK	7.144	7.144	0.000	1084977	20.0000	20.3
10 BENZALDEHYDE	8.164	8.164	0.000	766094	20.0000	20.2
11 ISO-PENTANAL	9.257	9.257	0.000	914971	20.0000	20.4
12 PENTANAL	9.944	9.944	0.000	899656	20.0000	20.4
14 O-TOLUALDEHYDE	11.017	11.017	0.000	725384	20.0000	20.0
15 M,P-TOLUALDEHYDE	11.324	11.324	0.000	1206012	40.0000	39.6
16 HEXANAL	12.850	12.850	0.000	789496	20.0000	19.6
17 2,5-Dimethylbenzaldehyde	13.444	13.444	0.000	564585	20.0000	19.0

Data File: /chem/hp1cf.i/14Sep2010.b/f0914031.d

Date : 15-SEP-2010 00:25

Client ID: 8

Sample Info: J1838-34-20.0:8

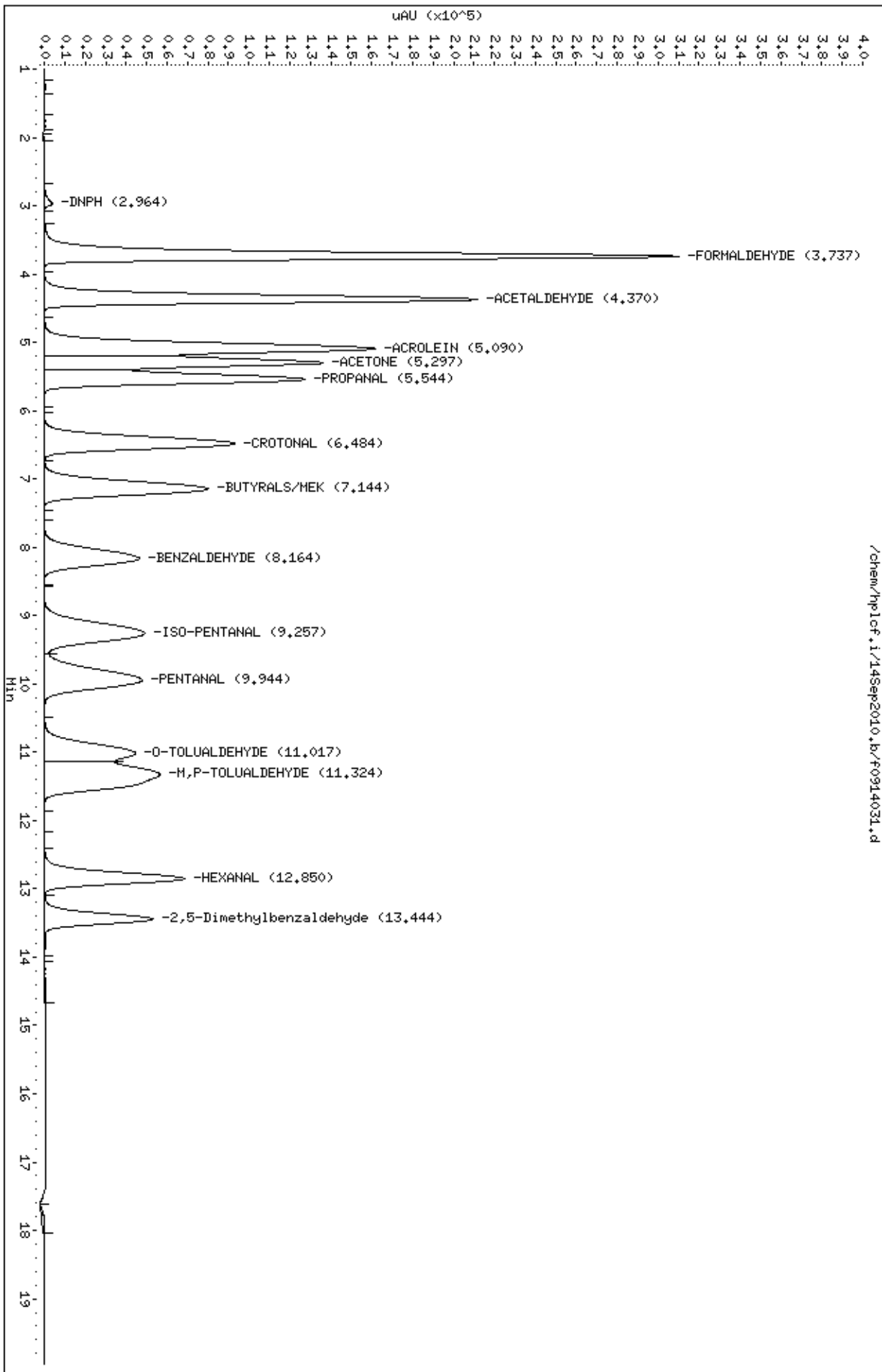
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914030.d
Lab Smp Id: 1838-34-20.0 Client Smp ID: 8
Inj Date : 15-SEP-2010 00:04
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-20.0;8
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 15-SEP-2010 00:04 Cal File: f0914030.d
Als bottle: 1 Calibration Sample, Level: 8
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.740	3.740	0.000	2254897	20.0000	21.0
3 ACETALDEHYDE	4.367	4.367	0.000	1716022	20.0000	20.4
4 ACRROLEIN	5.087	5.087	0.000	1624451	20.0000	20.3
5 ACETONE	5.294	5.294	0.000	1258629	20.0000	20.4
6 PROPANAL	5.540	5.540	0.000	1254219	20.0000	20.2
7 CROTONAL	6.481	6.481	0.000	1134607	20.0000	20.1
9 BUTYRALS/MEK	7.141	7.141	0.000	1085317	20.0000	20.3
10 BENZALDEHYDE	8.161	8.161	0.000	766359	20.0000	20.3
11 ISO-PENTANAL	9.254	9.254	0.000	915854	20.0000	20.4
12 PENTANAL	9.941	9.941	0.000	899974	20.0000	20.4
14 O-TOLUALDEHYDE	11.014	11.014	0.000	724371	20.0000	20.0
15 M,P-TOLUALDEHYDE	11.327	11.327	0.000	1208247	40.0000	39.6
16 HEXANAL	12.854	12.854	0.000	790179	20.0000	19.6
17 2,5-Dimethylbenzaldehyde	13.447	13.447	0.000	564439	20.0000	18.9

Data File: /chem/hp1cf.i/14Sep2010.b/f0914030.d

Date : 15-SEP-2010 00:04

Client ID: 8

Sample Info: J1838-34-20.0:8

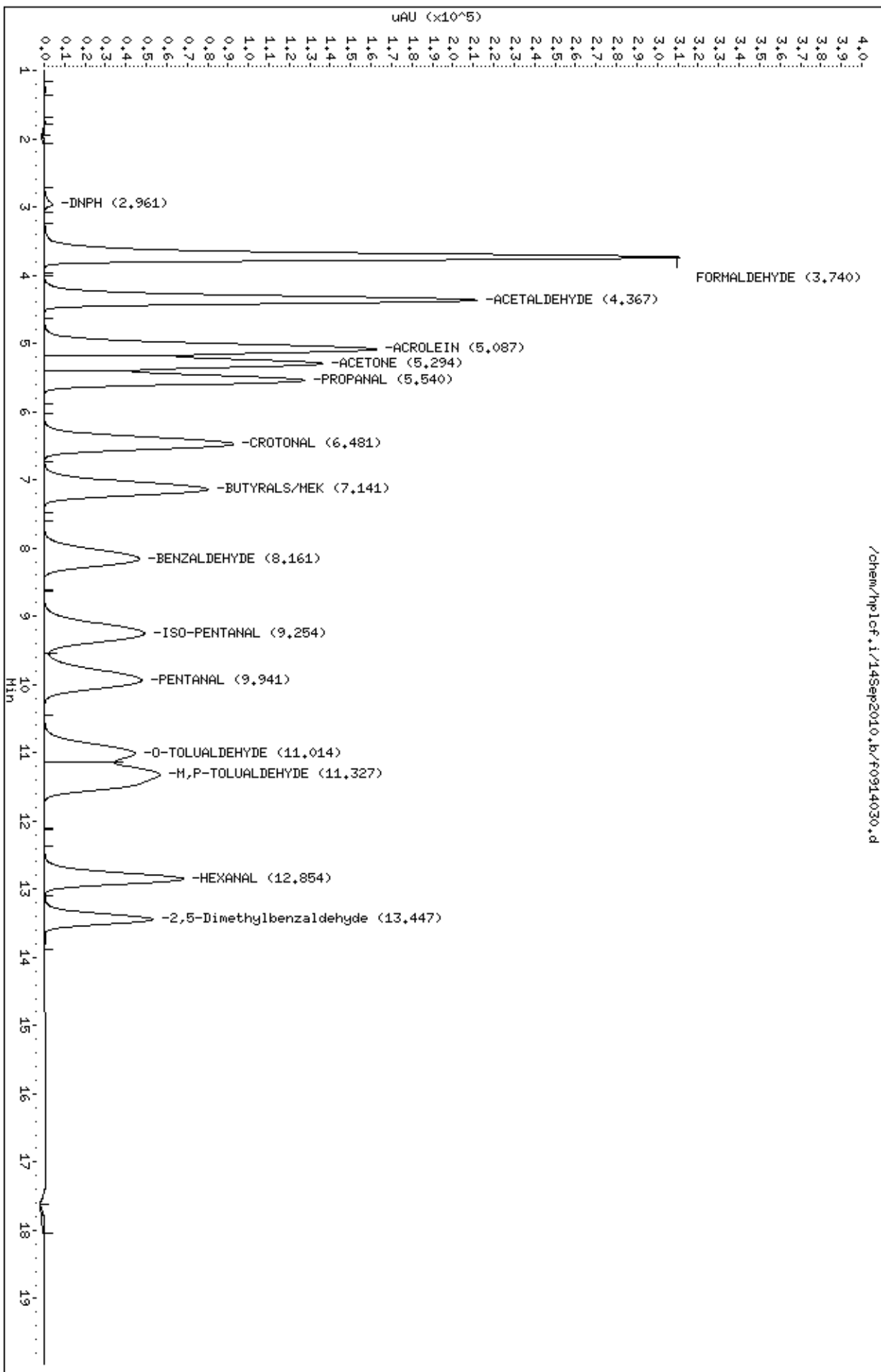
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914035.d
Lab Smp Id: 1838-34-50.0 Client Smp ID: 9
Inj Date : 15-SEP-2010 01:49
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-50.0;9
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 15-SEP-2010 01:49 Cal File: f0914035.d
Als bottle: 1 Calibration Sample, Level: 9
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.744	3.744	0.000	5735655	50.0000	52.9
3 ACETALDEHYDE	4.378	4.378	0.000	4320919	50.0000	51.1
4 ACRROLEIN	5.104	5.104	0.000	4122815	50.0000	51.3
5 ACETONE	5.311	5.311	0.000	3151407	50.0000	50.9
6 PROPANAL	5.558	5.558	0.000	3149158	50.0000	50.6
7 CROTONAL	6.498	6.498	0.000	2859994	50.0000	50.6
9 BUTYRALS/MEK	7.171	7.171	0.000	2736122	50.0000	51.0
10 BENZALDEHYDE	8.191	8.191	0.000	1932493	50.0000	50.9
11 ISO-PENTANAL	9.298	9.298	0.000	2308391	50.0000	51.2
12 PENTANAL	9.984	9.984	0.000	2269288	50.0000	51.2
14 O-TOLUALDEHYDE	11.058	11.058	0.000	1813745	50.0000	50.0
15 M,P-TOLUALDEHYDE	11.364	11.364	0.000	3063405	100.000	100
16 HEXANAL	12.878	12.878	0.000	1989961	50.0000	49.6
17 2,5-Dimethylbenzaldehyde	13.478	13.478	0.000	1422785	50.0000	48.3

Data File: /chem/hplcf.i/14Sep2010.b/f0914035.d

Date : 15-SEP-2010 01:49

Client ID: 9

Sample Info: f1838-34-50.0:9

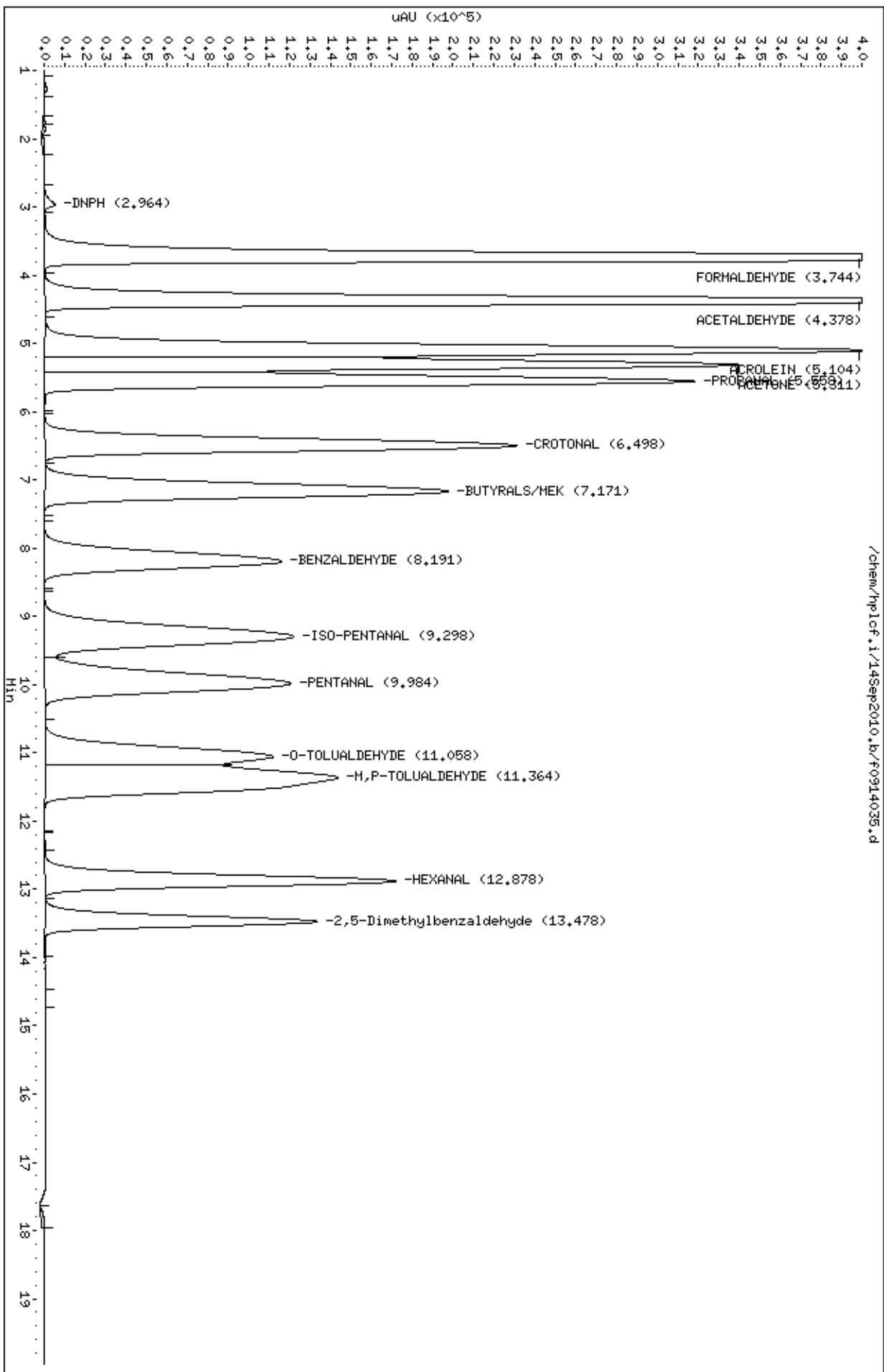
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914034.d
Lab Smp Id: 1838-34-50.0 Client Smp ID: 9
Inj Date : 15-SEP-2010 01:28
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-50.0;9
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 15-SEP-2010 01:28 Cal File: f0914034.d
Als bottle: 1 Calibration Sample, Level: 9
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.741	3.741	0.000	5712516	50.0000	52.8
3 ACETALDEHYDE	4.375	4.375	0.000	4302015	50.0000	50.9
4 ACRROLEIN	5.101	5.101	0.000	4126536	50.0000	51.4
5 ACETONE	5.308	5.308	0.000	3139620	50.0000	50.8
6 PROPANAL	5.555	5.555	0.000	3117295	50.0000	50.2
7 CROTONAL	6.495	6.495	0.000	2849453	50.0000	50.4
9 BUTYRALS/MEK	7.168	7.168	0.000	2731639	50.0000	50.9
10 BENZALDEHYDE	8.195	8.195	0.000	1923200	50.0000	50.7
11 ISO-PENTANAL	9.301	9.301	0.000	2299225	50.0000	51.1
12 PENTANAL	9.995	9.995	0.000	2257589	50.0000	51.0
14 O-TOLUALDEHYDE	11.075	11.075	0.000	1841301	50.0000	50.8
15 M,P-TOLUALDEHYDE	11.381	11.381	0.000	3009648	100.000	98.8
16 HEXANAL	12.888	12.888	0.000	1981435	50.0000	49.3
17 2,5-Dimethylbenzaldehyde	13.481	13.481	0.000	1415893	50.0000	48.0

Data File: /chem/hplcf.i/14Sep2010.b/f0914034.d

Date : 15-SEP-2010 01:28

Client ID: 9

Sample Info: J1838-34-50.0:9

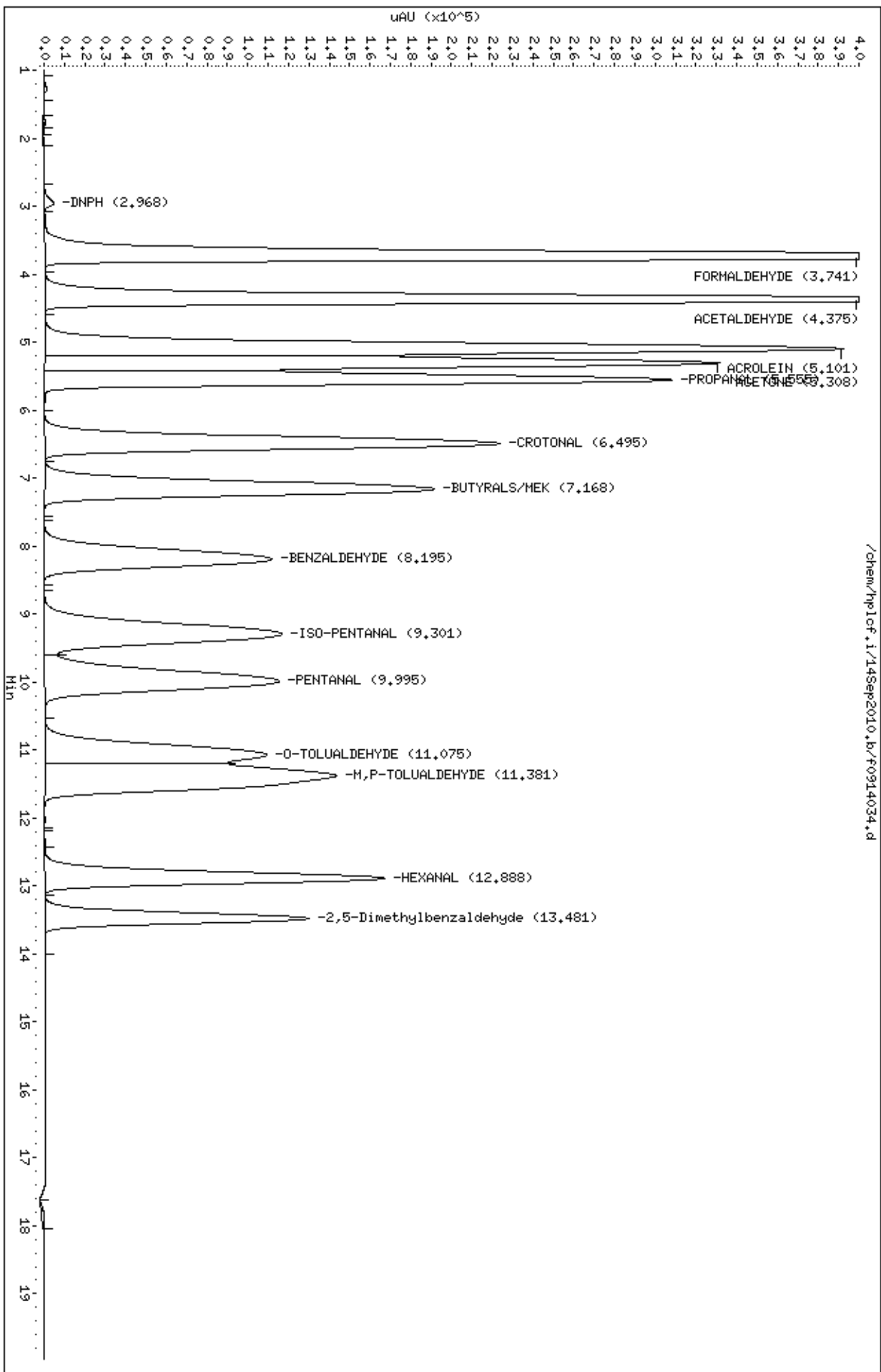
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914033.d
Lab Smp Id: 1838-34-50.0 Client Smp ID: 9
Inj Date : 15-SEP-2010 01:07
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-50.0;9
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 15-SEP-2010 01:07 Cal File: f0914033.d
Als bottle: 1 Calibration Sample, Level: 9
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (UG)	ON-COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.746	3.746	0.000	5728855	50.0000	53.0
3 ACETALDEHYDE	4.379	4.379	0.000	4315096	50.0000	51.1
4 ACRROLEIN	5.106	5.106	0.000	4140018	50.0000	51.6
5 ACETONE	5.312	5.312	0.000	3097047	50.0000	50.1
6 PROPANAL	5.553	5.553	0.000	3180515	50.0000	51.2
7 CROTONAL	6.499	6.499	0.000	2858190	50.0000	50.6
9 BUTYRALS/MEK	7.166	7.166	0.000	2734583	50.0000	51.0
10 BENZALDEHYDE	8.193	8.193	0.000	1931644	50.0000	51.0
11 ISO-PENTANAL	9.293	9.293	0.000	2306912	50.0000	51.3
12 PENTANAL	9.986	9.986	0.000	2269239	50.0000	51.3
14 O-TOLUALDEHYDE	11.053	11.053	0.000	1798798	50.0000	49.7
15 M,P-TOLUALDEHYDE	11.359	11.359	0.000	3071841	100.000	101
16 HEXANAL	12.879	12.879	0.000	1987958	50.0000	49.4
17 2,5-Dimethylbenzaldehyde	13.473	13.473	0.000	1421423	50.0000	48.0

Data File: /chem/hplcf.i/14Sep2010.b/f0914033.d

Date : 15-SEP-2010 01:07

Client ID: 9

Sample Info: J1838-34-50.0:9

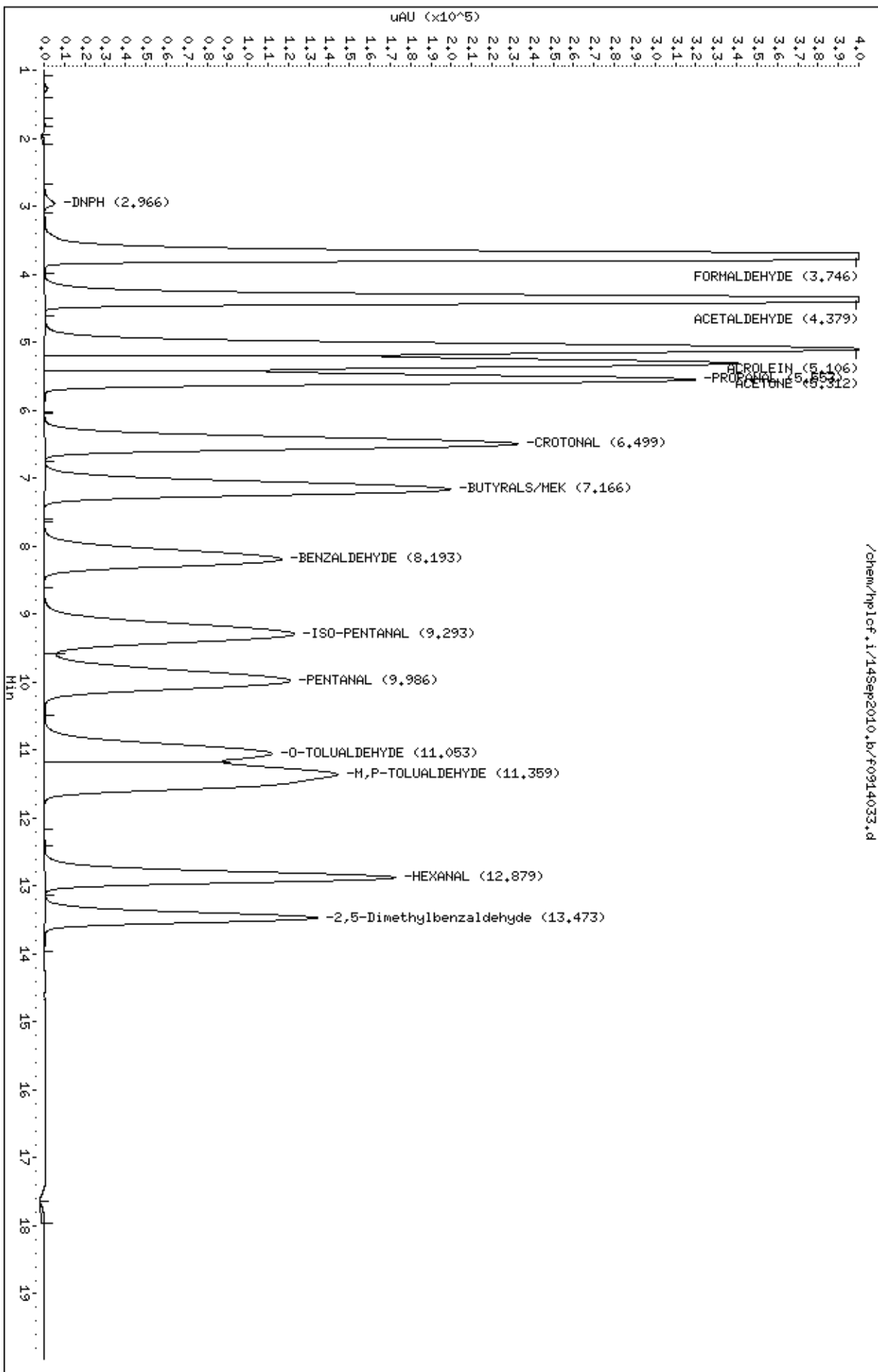
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914038.d
Lab Smp Id: 1838-34-75.0 Client Smp ID: 10
Inj Date : 15-SEP-2010 02:52
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-75.0;10
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 15-SEP-2010 02:52 Cal File: f0914038.d
Als bottle: 1 Calibration Sample, Level: 10
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL - AMT (UG)	ON - COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.740	3.740	0.000	8877830	75.0000	81.1(A)
3 ACETALDEHYDE	4.373	4.373	0.000	6766583	75.0000	79.4(A)
4 ACRROLEIN	5.093	5.093	0.000	6437988	75.0000	79.4(A)
5 ACETONE	5.300	5.300	0.000	4958940	75.0000	79.6(A)
6 PROPANAL	5.546	5.546	0.000	4945549	75.0000	78.9(A)
7 CROTONAL	6.480	6.480	0.000	4483386	75.0000	78.7(A)
9 BUTYRALS/MEK	7.146	7.146	0.000	4290345	75.0000	79.2(A)
10 BENZALDEHYDE	8.166	8.166	0.000	3033347	75.0000	79.3(A)
11 ISO-PENTANAL	9.260	9.260	0.000	3618335	75.0000	79.6(A)
12 PENTANAL	9.946	9.946	0.000	3558158	75.0000	79.6(A)
14 O-TOLUALDEHYDE	11.026	11.026	0.000	2834077	75.0000	77.7(A)
15 M,P-TOLUALDEHYDE	11.333	11.333	0.000	4808540	150.000	157(A)
16 HEXANAL	12.853	12.853	0.000	3118236	75.0000	77.3(A)
17 2,5-Dimethylbenzaldehyde	13.453	13.453	0.000	2231351	75.0000	75.6(A)

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.

Data File: /chem/hp1cf.i/14Sep2010.b/f0914038.d

Date : 15-SEP-2010 02:52

Client ID: 10

Sample Info: f1838-34-75.0:10

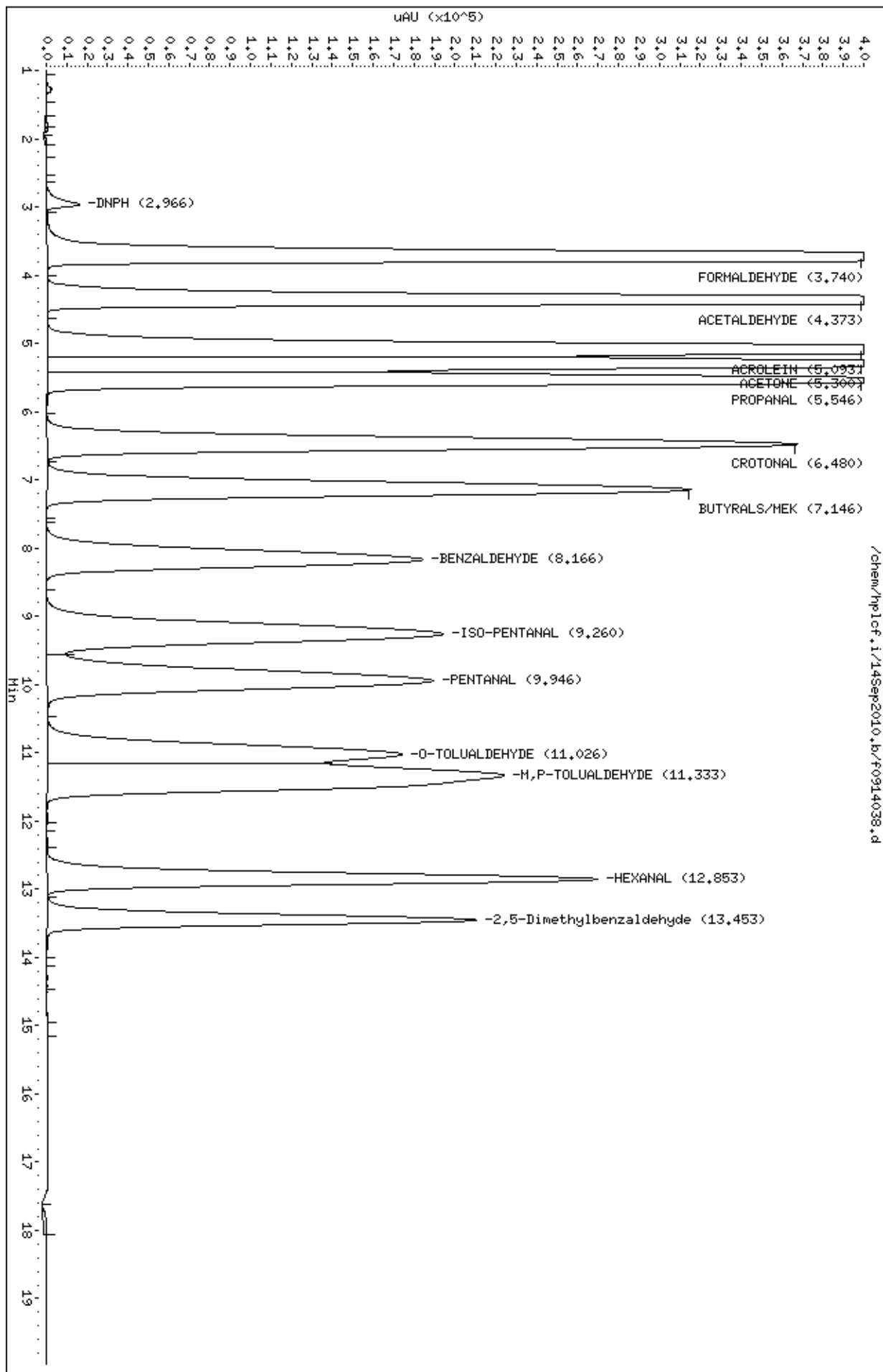
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914037.d
Lab Smp Id: 1838-34-75.0 Client Smp ID: 10
Inj Date : 15-SEP-2010 02:31
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-75.0;10
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 15-SEP-2010 02:31 Cal File: f0914037.d
Als bottle: 1 Calibration Sample, Level: 10
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL - AMT (UG)	ON - COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.743	3.743	0.000	8874471	75.0000	81.3(A)
3 ACETALDEHYDE	4.376	4.376	0.000	6764082	75.0000	79.6(A)
4 ACROLEIN	5.103	5.103	0.000	6468876	75.0000	79.9(A)
5 ACETONE	5.310	5.310	0.000	4877086	75.0000	78.5(A)
6 PROPANAL	5.550	5.550	0.000	4992133	75.0000	79.8(A)
7 CROTONAL	6.496	6.496	0.000	4481980	75.0000	78.9(A)
9 BUTYRALS/MEK	7.163	7.163	0.000	4289038	75.0000	79.4(A)
10 BENZALDEHYDE	8.183	8.183	0.000	3028322	75.0000	79.4(A)
11 ISO-PENTANAL	9.283	9.283	0.000	3619020	75.0000	79.8(A)
12 PENTANAL	9.970	9.970	0.000	3557001	75.0000	79.8(A)
14 O-TOLUALDEHYDE	11.043	11.043	0.000	2868309	75.0000	78.8(A)
15 M,P-TOLUALDEHYDE	11.356	11.356	0.000	4770723	150.000	156(A)
16 HEXANAL	12.870	12.870	0.000	3117337	75.0000	77.4(A)
17 2,5-Dimethylbenzaldehyde	13.470	13.470	0.000	2229525	75.0000	75.6(A)

QC Flag Legend

A - Target compound detected but, quantitated amount
exceeded maximum amount.

Data File: /chem/hp1cf.i/14Sep2010.b/f0914037.d

Date : 15-SEP-2010 02:31

Client ID: 10

Sample Info: f1838-34-75.0:10

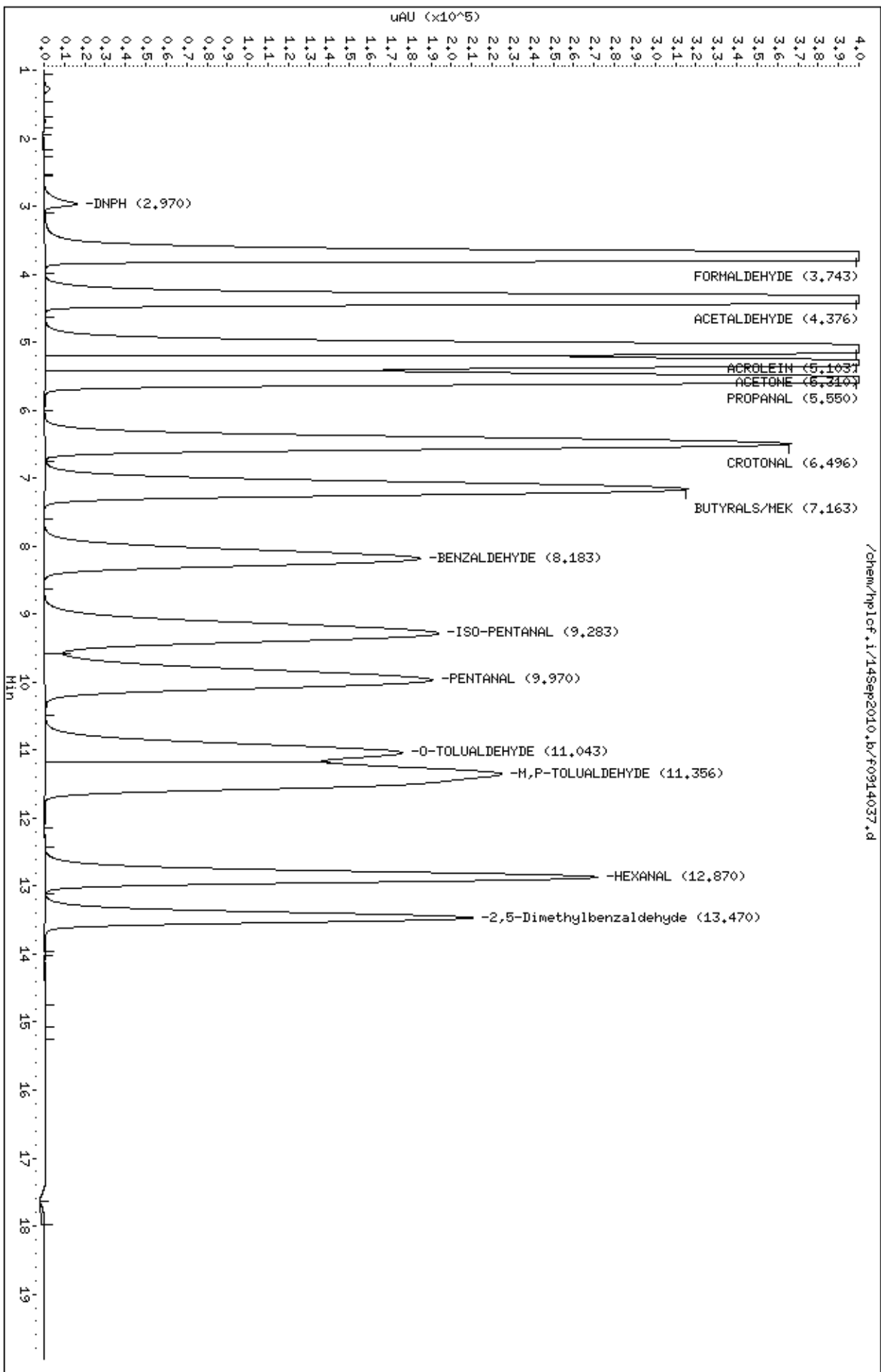
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/14Sep2010.b/f0914036.d
Lab Smp Id: 1838-34-75.0 Client Smp ID: 10
Inj Date : 15-SEP-2010 02:10
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-34-75.0;10
Misc Info :
Comment :
Method : /chem/hplcf.i/14Sep2010.b/F10D0914.m
Meth Date : 15-Sep-2010 09:40 cleaf Quant Type: ESTD
Cal Date : 15-SEP-2010 02:10 Cal File: f0914036.d
Als bottle: 1 Calibration Sample, Level: 10
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable
Vs 5.00000 Sample Final Volume (mL)
Vo 5.00000 Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL - AMT (UG)	ON - COL (UG)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.758	3.758	0.000	8872541	75.0000	81.5(A)
3 ACETALDEHYDE	4.391	4.391	0.000	6752657	75.0000	79.7(A)
4 ACRROLEIN	5.125	5.125	0.000	6480467	75.0000	80.3(A)
5 ACETONE	5.332	5.332	0.000	4840692	75.0000	78.0(A)
6 PROPANAL	5.578	5.578	0.000	4977690	75.0000	79.8(A)
7 CROTONAL	6.525	6.525	0.000	4481850	75.0000	79.1(A)
9 BUTYRALS/MEK	7.198	7.198	0.000	4287442	75.0000	79.6(A)
10 BENZALDEHYDE	8.238	8.238	0.000	3023220	75.0000	79.4(A)
11 ISO-PENTANAL	9.345	9.345	0.000	3616516	75.0000	80.0(A)
12 PENTANAL	10.038	10.038	0.000	3551905	75.0000	79.9(A)
14 O-TOLUALDEHYDE	11.098	11.098	0.000	2865895	75.0000	78.9(A)
15 M,P-TOLUALDEHYDE	11.405	11.405	0.000	4759434	150.000	156(A)
16 HEXANAL	12.912	12.912	0.000	3113906	75.0000	77.4(A)
17 2,5-Dimethylbenzaldehyde	13.505	13.505	0.000	2224010	75.0000	75.4(A)

QC Flag Legend

A - Target compound detected but, quantitated amount
exceeded maximum amount.

Data File: /chem/hp1cf.i/14Sep2010.b/f0914036.d

Date : 15-SEP-2010 02:10

Client ID: 10

Sample Info: f1838-34-75.0:10

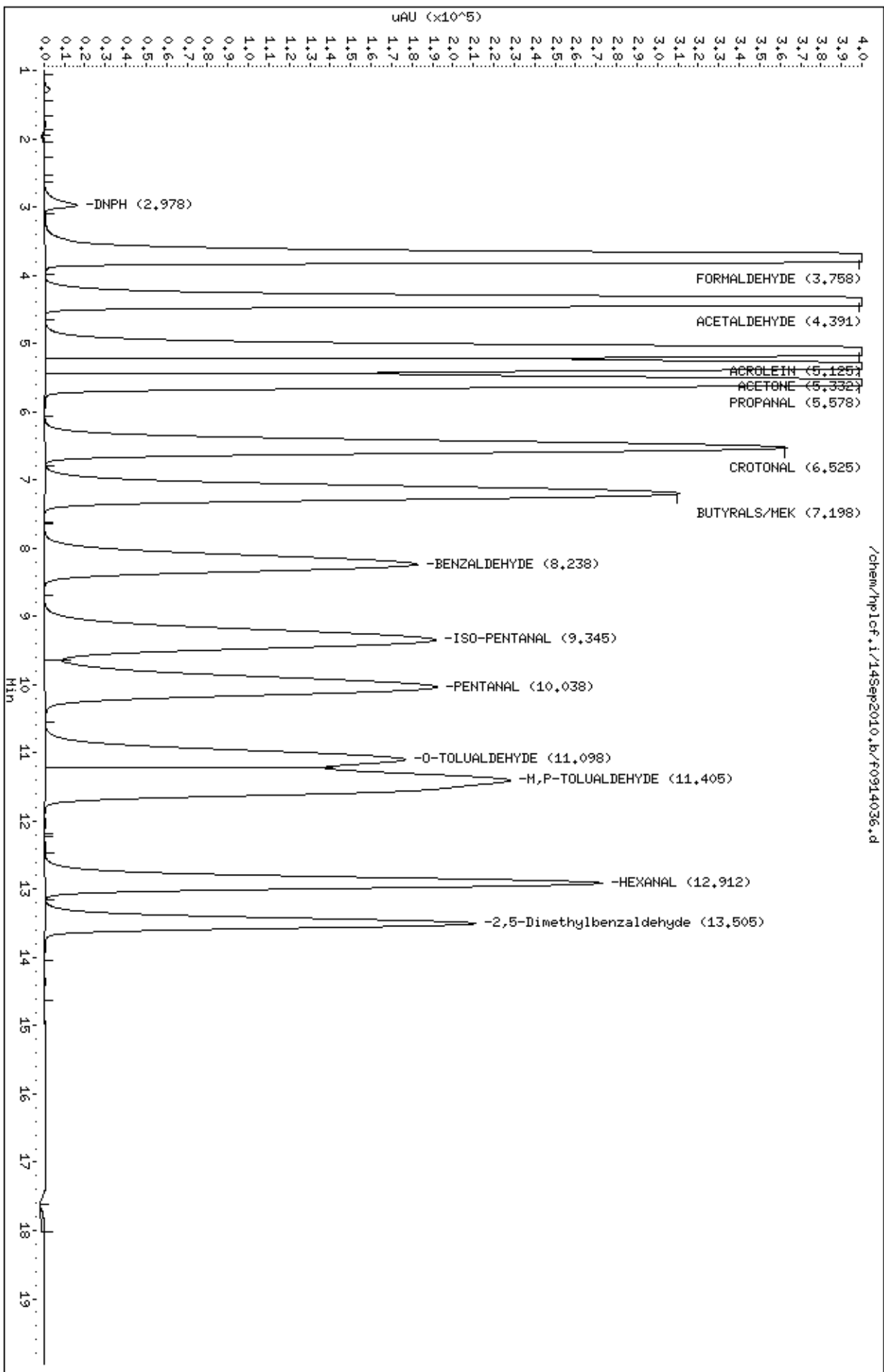
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

CONTINUING CALIBRATION COMPOUNDS

Instrument ID: hplcf.i Injection Date: 09-NOV-2010 09:19
Lab File ID: f1109003.d Init. Cal. Date(s): 14-SEP-2010 15-SEP-2010
Analysis Type: AIR Init. Cal. Times: 16:46 02:52
Lab Sample ID: 1838-41-10 Quant Type: ESTD
Method: /chem/hplcf.i/09Nov2010.b/F10D0914.m

COMPOUND	RRF / AMOUNT	RF10	MIN RRF	%D / %DRIFT	MAX %D / %DRIFT	CURVE TYPE
1 DNPH	++++	387	0.100	++++	10.00000	Averaged
2 FORMALDEHYDE	109474	115283	0.010	-5.30611	10.00000	Averaged
3 ACETALDEHYDE	85159	86425	0.010	-1.48656	10.00000	Averaged
4 ACROLEIN	81132	84443	0.010	-4.08213	10.00000	Averaged
5 ACETONE	62324	61761	0.010	0.90439	10.00000	Averaged
6 PROPANAL	62684	61536	0.010	1.83238	10.00000	Averaged
7 CROTONAL	56944	57195	0.010	-0.43989	10.00000	Averaged
9 BUTYRALS/MEK	54133	53871	0.010	0.48397	10.00000	Averaged
10 BENZALDEHYDE	38259	37711	0.010	1.43245	10.00000	Averaged
11 ISO-PENTANAL	45456	45603	0.010	-0.32286	10.00000	Averaged
12 PENTANAL	44700	44397	0.010	0.67941	10.00000	Averaged
14 O-TOLUALDEHYDE	36466	36812	0.010	-0.94906	10.00000	Averaged
15 M,P-TOLUALDEHYDE	30634	29926	0.010	2.30973	10.00000	Averaged
16 HEXANAL	40334	39843	0.010	1.21704	10.00000	Averaged
17 2,5-Dimethylbenzaldehyde	29495	28666	0.010	2.81042	10.00000	Averaged

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109003.d
Lab Smp Id: 1838-41-10 Client Smp ID: CCV
Inj Date : 09-NOV-2010 09:19
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-41-10;CCV
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 10:28 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1 Continuing Calibration Sample
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	5.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	AMOUNTS	
					CAL-AMT (ng)	ON-COL (ng)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.731	3.731	0.000	1152831	10.0000	10.5
3 ACETALDEHYDE	4.344	4.344	0.000	864245	10.0000	10.1
4 ACROLEIN	5.051	5.051	0.000	844434	10.0000	10.4
5 ACETONE	5.250	5.250	0.000	617606	10.0000	9.91
6 PROPANAL	5.497	5.497	0.000	615355	10.0000	9.82
7 CROTONAL	6.417	6.417	0.000	571946	10.0000	10.0
9 BUTYRALS/MEK	7.064	7.064	0.000	538712	10.0000	9.95
10 BENZALDEHYDE	8.064	8.064	0.000	377108	10.0000	9.86
11 ISO-PENTANAL	9.104	9.104	0.000	456028	10.0000	10.0
12 PENTANAL	9.764	9.764	0.000	443965	10.0000	9.93
14 O-TOLUALDEHYDE	10.877	10.877	0.000	368122	10.0000	10.1
15 M,P-TOLUALDEHYDE	11.184	11.184	0.000	598524	20.0000	19.5
16 HEXANAL	12.717	12.717	0.000	398431	10.0000	9.88
17 2,5-Dimethylbenzaldehyde	13.324	13.324	0.000	286660	10.0000	9.72

Data File: /chem/hplcf.i/09Nov2010.b/f1109003.d

Date : 09-NOV-2010 09:19

Client ID: CCV

Sample Info: f1838-41-10;CCV

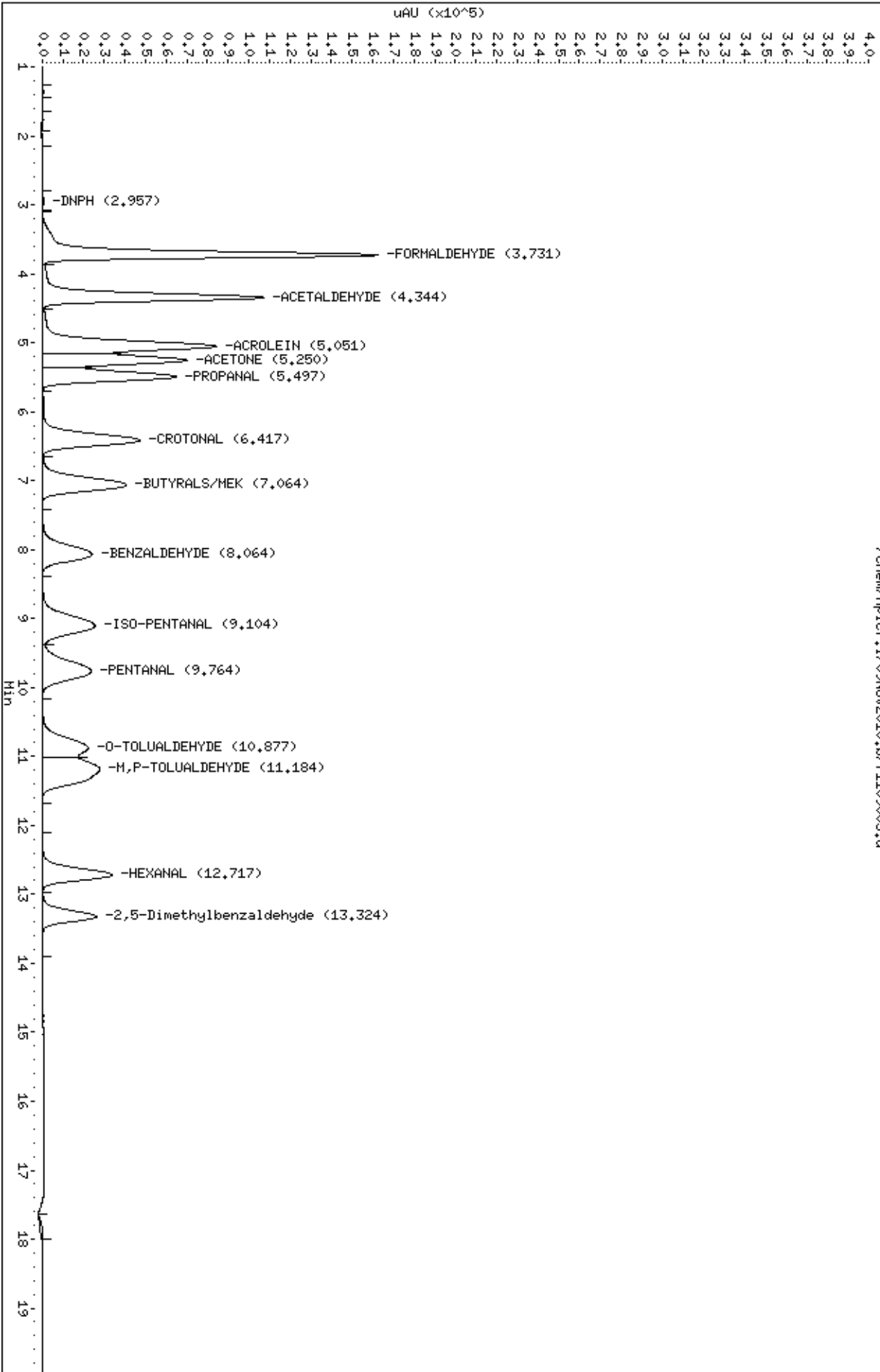
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: CRL

Column diameter: 2.00



```

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109004.d
Lab Smp Id: CCB                               Client Smp ID: LAB BLANK
Inj Date   : 09-NOV-2010 09:40
Operator   : CRL                               Inst ID: hplcf.i
Smp Info   : ;CCB;LAB BLANK
Misc Info  :
Comment    :
Method     : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date  : 09-Nov-2010 10:28 cleaf          Quant Type: ESTD
Cal Date   : 14-SEP-2010 18:30               Cal File: f0914014.d
Als bottle: 1
Dil Factor: 1.00000
Integrator: Falcon                            Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

```

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	5.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	CONCENTRATIONS	
							ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====			=====	=====	
1 DNPH		Compound	Not	Detected.				
2 FORMALDEHYDE		Compound	Not	Detected.				
3 ACETALDEHYDE		Compound	Not	Detected.				
4 ACROLEIN		Compound	Not	Detected.				
5 ACETONE		Compound	Not	Detected.				
6 PROPANAL		Compound	Not	Detected.				
7 CROTONAL		Compound	Not	Detected.				
9 BUTYRALS/MEK		Compound	Not	Detected.				
10 BENZALDEHYDE		Compound	Not	Detected.				
11 ISO-PENTANAL		Compound	Not	Detected.				
12 PENTANAL		Compound	Not	Detected.				
14 O-TOLUALDEHYDE		Compound	Not	Detected.				
15 M,P-TOLUALDEHYDE		Compound	Not	Detected.				
16 HEXANAL		Compound	Not	Detected.				

Compounds	RT	CONCENTRATIONS			RESPONSE	ON-COLUMN		FINAL
		EXP	RT	DLT		(	ng)	(
=====	==	=====	=====		=====	=====		=====
17 2,5-Dimethylbenzaldehyde					Compound Not Detected.			

Data File: /chem/hp1cf.i/09Nov2010.b/f1109004.d

Page 1

Date : 09-NOV-2010 09:40

Client ID: LAB BLANK

Instrument: hp1cf.i

Sample Info: JCCB;LAB BLANK

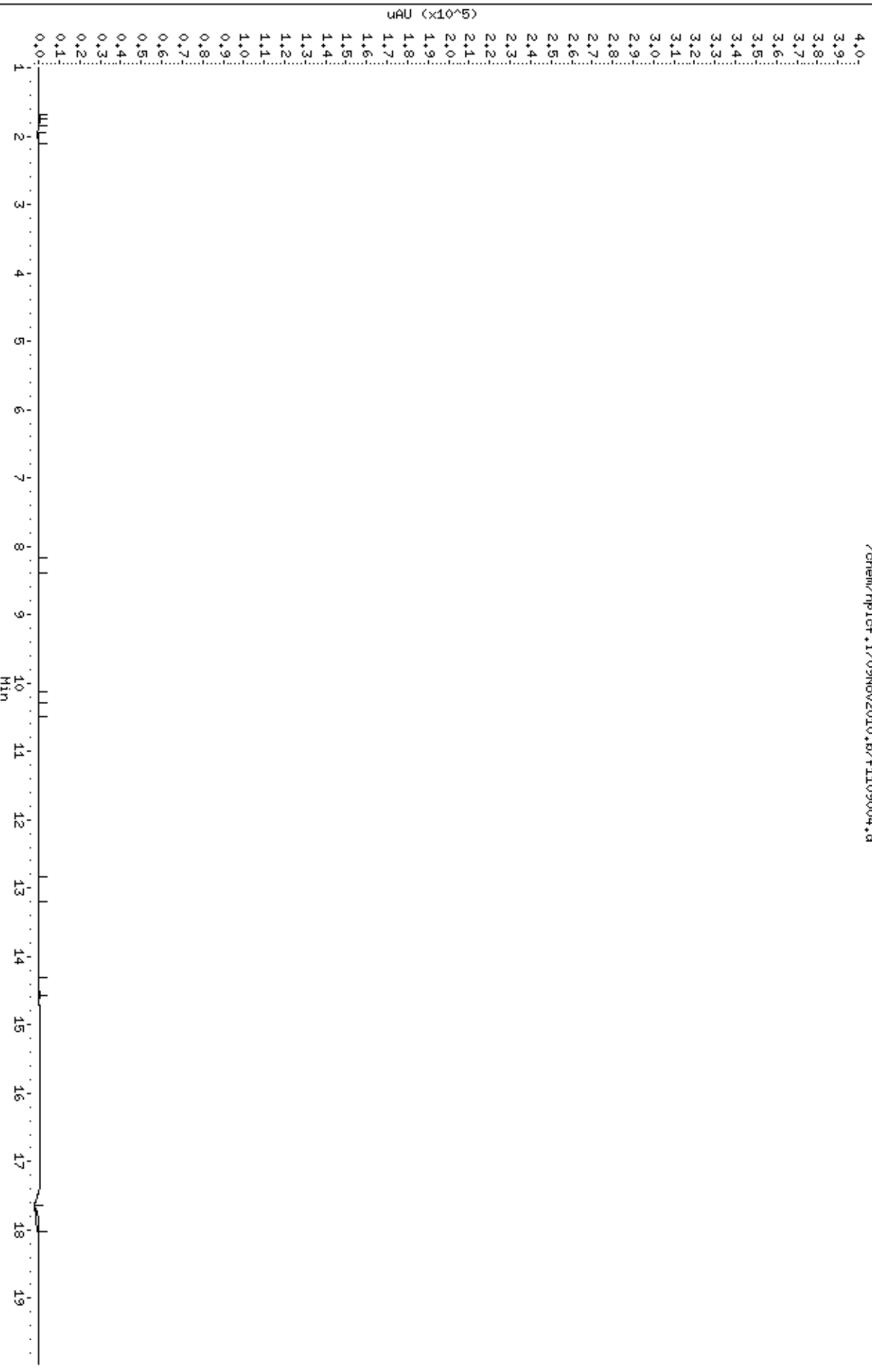
Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00

/chem/hp1cf.i/09Nov2010.b/f1109004.d



Air Toxics Ltd.

TARGET COMPOUNDS

Client Name:
Lab Smp Id: 1838-41-10
Sample Location:
Sample Date:
Sample Matrix: AIR
Analysis Type: OTHER
Data Type: LC DATA
Misc Info:

Client SDG: 09Nov2010
Client Smp ID: CCV
Sample Point:
Date Received:
Quant Type: ESTD
Level: MED
Operator: CRL

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/KG) ug	Q
119-26-6-----	DNPH	0.00	
50-00-0-----	FORMALDEHYDE	10.5	
75-07-0-----	ACETALDEHYDE	10.1	
107-02-8-----	ACROLEIN	10.5	
67-64-1-----	ACETONE	9.70	
123-38-6-----	PROPANAL	9.80	
123-73-9-----	CROTONAL	9.99	
9999-9999-007--	BUTYRALS/MEK	9.91	
100-52-7-----	BENZALDEHYDE	9.85	
590-86-3-----	ISO-PENTANAL	10.0	
110-62-3-----	PENTANAL	9.89	
529-20-4-----	O-TOLUALDEHYDE	10.1	
620-23-51-----	M, P-TOLUALDEHYDE	19.6	
66-25-1-----	HEXANAL	9.91	
5779-94-2-----	2,5-Dimethylbenzaldehyde	9.75	
=====	=====	=====	=====

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109015.d
Lab Smp Id: 1838-41-10 Client Smp ID: CCV
Inj Date : 09-NOV-2010 13:30
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-41-10;CCV
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1 QC Sample: METHSPIKE
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	5.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds					CONCENTRATIONS	
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.737	3.731	0.006	1152228	10.5251	10.5
3 ACETALDEHYDE	4.357	4.344	0.013	862012	10.1224	10.1
4 ACRROLEIN	5.071	5.051	0.020	853207	10.5164	10.5
5 ACETONE	5.271	5.250	0.021	604903	9.70575	9.70
6 PROPANAL	5.511	5.497	0.014	614360	9.80089	9.80
7 CROTONAL	6.431	6.417	0.014	568720	9.98734	9.99
9 BUTYRALS/MEK	7.084	7.064	0.020	536589	9.91238	9.91
10 BENZALDEHYDE	8.084	8.064	0.020	377013	9.85427	9.85
11 ISO-PENTANAL	9.124	9.104	0.020	456476	10.0421	10.0
12 PENTANAL	9.791	9.764	0.027	441956	9.88712	9.89
14 O-TOLUALDEHYDE	10.897	10.877	0.020	367073	10.0662	10.1
15 M,P-TOLUALDEHYDE	11.211	11.184	0.027	600255	19.5946	19.6
16 HEXANAL	12.737	12.717	0.020	399588	9.90697	9.91
17 2,5-Dimethylbenzaldehyde	13.351	13.324	0.027	287608	9.75111	9.75

Air Toxics Ltd.

RECOVERY REPORT

Client Name: Client SDG: 09Nov2010
Sample Matrix: LIQUID Fraction: OTHER
Lab Smp Id: 1838-41-10 Client Smp ID: CCV
Level: MED Operator: CRL
Data Type: LC DATA SampleType: METHSPIKE
SpikeList File: 10ug.spk Quant Type: ESTD
Sublist File: all.sub
Method File: /chem/hplcf.i/09Nov2010.b/F10D0914.m
Misc Info:

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	10.0	10.5	105.25	90-110
3 ACETALDEHYDE	10.0	10.1	101.22	90-110
4 ACROLEIN	10.0	10.5	105.16	90-110
5 ACETONE	10.0	9.70	97.06	90-110
6 PROPANAL	10.0	9.80	98.01	90-110
7 CROTONAL	10.0	9.99	99.87	90-110
9 BUTYRALS/MEK	10.0	9.91	99.12	90-110
10 BENZALDEHYDE	10.0	9.85	98.54	90-110
11 ISO-PENTANAL	10.0	10.0	100.42	90-110
12 PENTANAL	10.0	9.89	98.87	90-110
14 O-TOLUALDEHYDE	10.0	10.1	100.66	90-110
15 M,P-TOLUALDEHYDE	20.0	19.6	97.97	90-110
16 HEXANAL	10.0	9.91	99.07	90-110
17 2,5-Dimethylbenzal	10.0	9.75	97.51	90-110

Data File: /chem/hp1cf.i/09Nov2010.b/f1109015.d

Date : 09-NOV-2010 13:30

Client ID: CCV

Sample Info: #1838-41-10;CCV

Purge Volume: 5.0

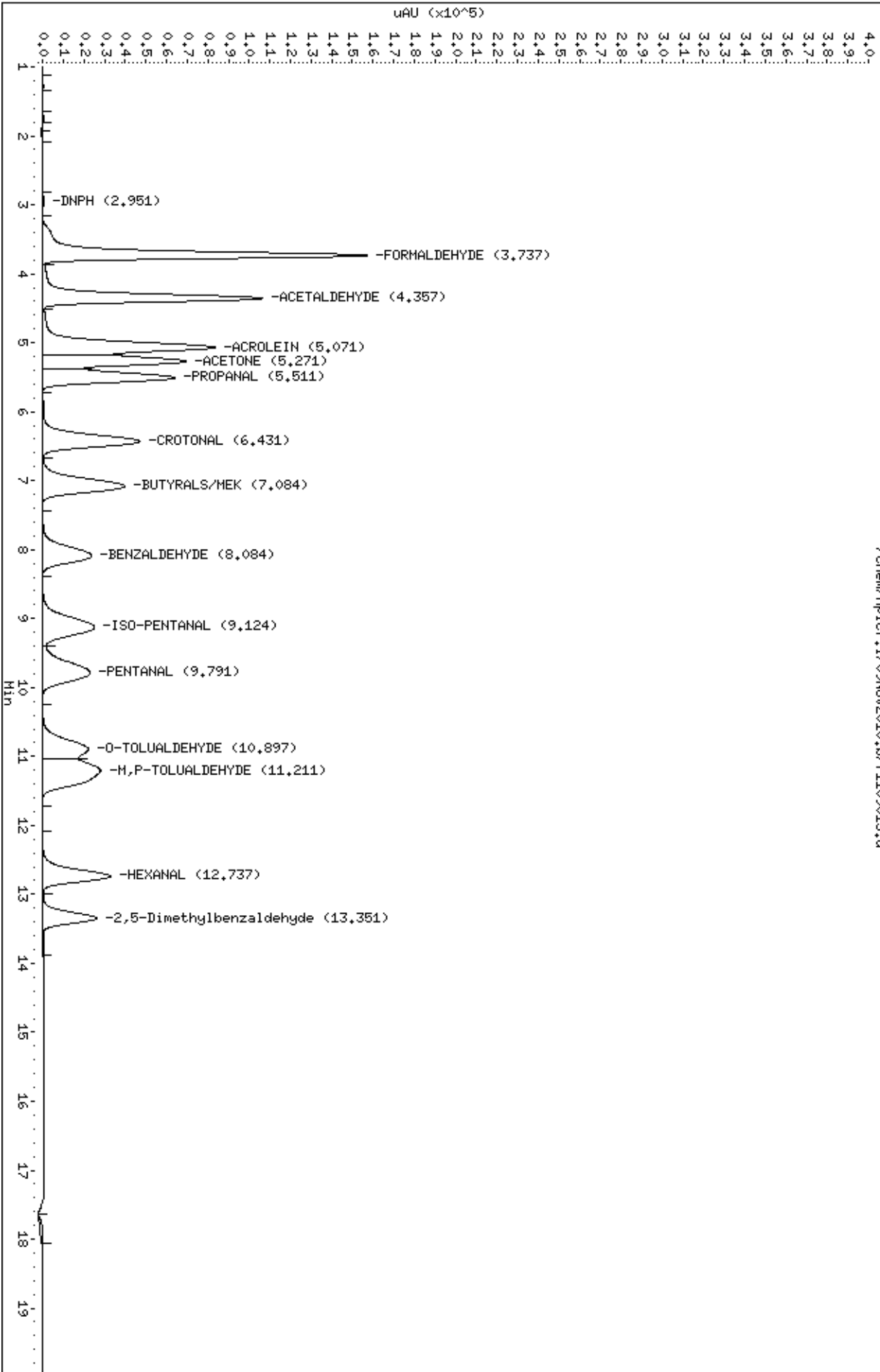
Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00

/chem/hp1cf.i/09Nov2010.b/f1109015.d



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109016.d
Lab Smp Id: CCB Client Smp ID: LAB BLANK
Inj Date : 09-NOV-2010 13:51
Operator : CRL Inst ID: hplcf.i
Smp Info : ;CCB;LAB BLANK
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	5.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	CONCENTRATIONS	
							ON-COLUMN	FINAL
							(ng)	(ug)
1 DNPH						Compound Not Detected.		
2 FORMALDEHYDE						Compound Not Detected.		
3 ACETALDEHYDE						Compound Not Detected.		
4 ACROLEIN						Compound Not Detected.		
5 ACETONE						Compound Not Detected.		
6 PROPANAL						Compound Not Detected.		
7 CROTONAL						Compound Not Detected.		
9 BUTYRALS/MEK						Compound Not Detected.		
10 BENZALDEHYDE						Compound Not Detected.		
11 ISO-PENTANAL						Compound Not Detected.		
12 PENTANAL						Compound Not Detected.		
14 O-TOLUALDEHYDE						Compound Not Detected.		
15 M,P-TOLUALDEHYDE						Compound Not Detected.		
16 HEXANAL						Compound Not Detected.		

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	CONCENTRATIONS	
							ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====			=====	=====	
17 2,5-Dimethylbenzaldehyde		Compound Not Detected.						

Data File: /chem/hplcf.i/09Nov2010.b/f1109016.d

Date : 09-NOV-2010 13:51

Client ID: LAB BLANK

Sample Info: JCCB;LAB BLANK

Purge Volume: 5.0

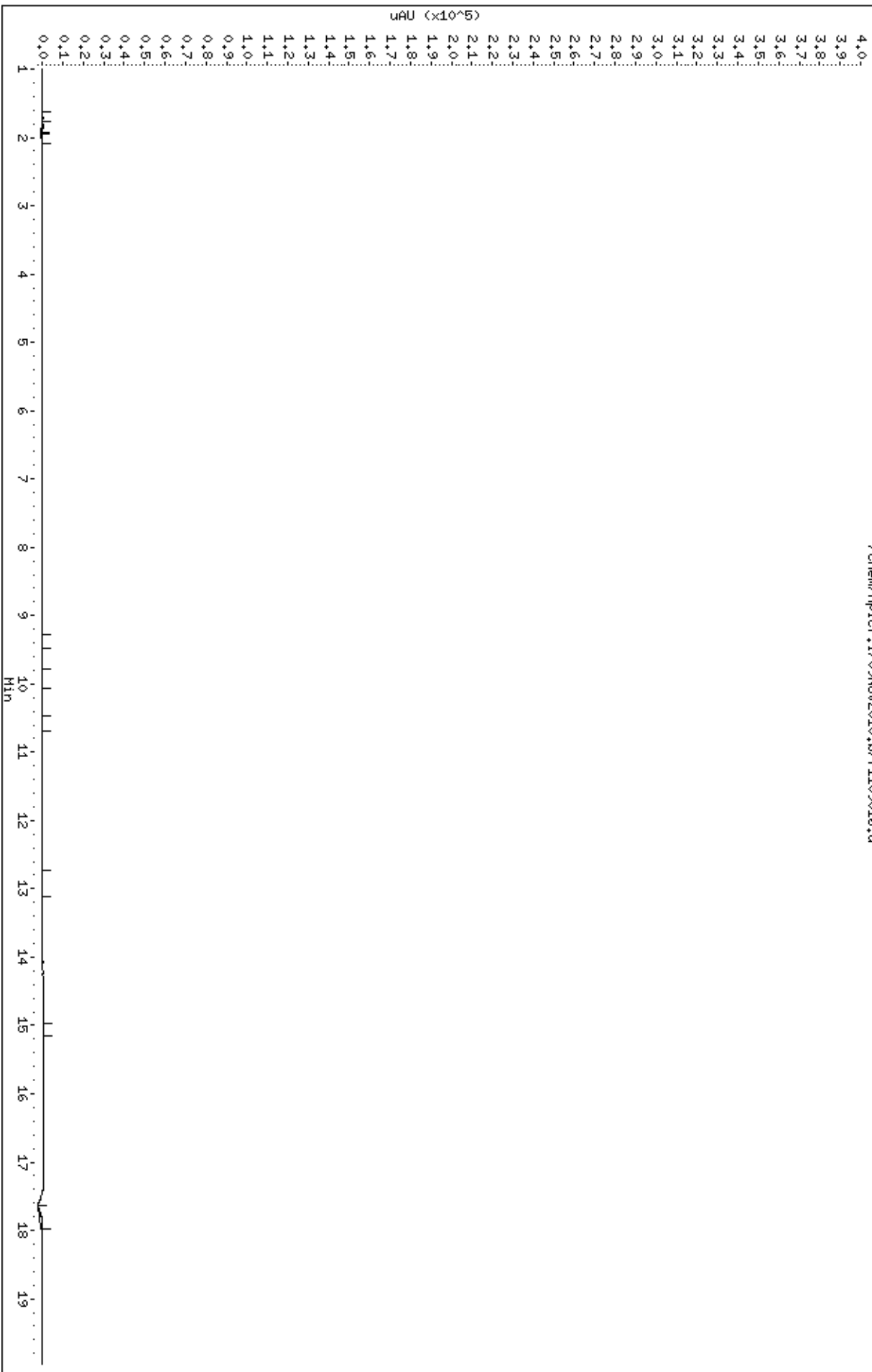
Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: CRL

Column diameter: 2.00

/chem/hplcf.i/09Nov2010.b/f1109016.d



Air Toxics Ltd.

TARGET COMPOUNDS

Client Name:	Client SDG: 09Nov2010
Lab Smp Id: 1838-41-10	Client Smp ID: CCV
Sample Location:	Sample Point:
Sample Date:	Date Received:
Sample Matrix: AIR	Quant Type: ESTD
Analysis Type: OTHER	Level: MED
Data Type: LC DATA	Operator: CRL
Misc Info:	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/KG) ug	Q
119-26-6-----	DNPH	0.00	
50-00-0-----	FORMALDEHYDE	10.4	
75-07-0-----	ACETALDEHYDE	10.0	
107-02-8-----	ACROLEIN	10.4	
67-64-1-----	ACETONE	9.71	
123-38-6-----	PROPANAL	9.59	
123-73-9-----	CROTONAL	9.85	
9999-9999-007--	BUTYRALS/MEK	9.82	
100-52-7-----	BENZALDEHYDE	9.80	
590-86-3-----	ISO-PENTANAL	9.97	
110-62-3-----	PENTANAL	9.81	
529-20-4-----	O-TOLUALDEHYDE	9.82	
620-23-51-----	M, P-TOLUALDEHYDE	19.5	
66-25-1-----	HEXANAL	9.86	
5779-94-2-----	2,5-Dimethylbenzaldehyde	10.1	
=====	=====	=====	=====

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
 Data file : /chem/hplcf.i/09Nov2010.b/f1109027.d
 Lab Smp Id: 1838-41-10 Client Smp ID: CCV
 Inj Date : 09-NOV-2010 17:40
 Operator : CRL Inst ID: hplcf.i
 Smp Info : ;1838-41-10;CCV
 Misc Info :
 Comment :
 Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
 Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
 Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
 Als bottle: 1 QC Sample: METHSPIKE
 Dil Factor: 1.00000
 Integrator: Falcon Compound Sublist: all.sub
 Target Version: 3.50
 Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	5.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds					CONCENTRATIONS	
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.691	3.731	-0.040	1144006	10.4500	10.4
3 ACETALDEHYDE	4.290	4.344	-0.054	853378	10.0210	10.0
4 ACRROLEIN	4.971	5.051	-0.080	841901	10.3770	10.4
5 ACETONE	5.170	5.250	-0.080	605355	9.71300	9.71
6 PROPANAL	5.404	5.497	-0.093	601439	9.59476	9.59
7 CROTONAL	6.284	6.417	-0.133	560739	9.84719	9.85
9 BUTYRALS/MEK	6.910	7.064	-0.154	531660	9.82132	9.82
10 BENZALDEHYDE	7.864	8.064	-0.200	374782	9.79595	9.80
11 ISO-PENTANAL	8.870	9.104	-0.234	453393	9.97432	9.97
12 PENTANAL	9.497	9.764	-0.267	438421	9.80804	9.81
14 O-TOLUALDEHYDE	10.624	10.877	-0.253	358212	9.82314	9.82
15 M,P-TOLUALDEHYDE	10.944	11.184	-0.240	598851	19.5487	19.5
16 HEXANAL	12.544	12.717	-0.173	397905	9.86525	9.86
17 2,5-Dimethylbenzaldehyde	13.164	13.324	-0.160	297744	10.0947	10.1

Air Toxics Ltd.

RECOVERY REPORT

Client Name: Client SDG: 09Nov2010
Sample Matrix: LIQUID Fraction: OTHER
Lab Smp Id: 1838-41-10 Client Smp ID: CCV
Level: MED Operator: CRL
Data Type: LC DATA SampleType: METHSPIKE
SpikeList File: 10ug.spk Quant Type: ESTD
Sublist File: all.sub
Method File: /chem/hplcf.i/09Nov2010.b/F10D0914.m
Misc Info:

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	10.0	10.4	104.50	90-110
3 ACETALDEHYDE	10.0	10.0	100.21	90-110
4 ACROLEIN	10.0	10.4	103.77	90-110
5 ACETONE	10.0	9.71	97.13	90-110
6 PROPANAL	10.0	9.59	95.95	90-110
7 CROTONAL	10.0	9.85	98.47	90-110
9 BUTYRALS/MEK	10.0	9.82	98.21	90-110
10 BENZALDEHYDE	10.0	9.80	97.96	90-110
11 ISO-PENTANAL	10.0	9.97	99.74	90-110
12 PENTANAL	10.0	9.81	98.08	90-110
14 O-TOLUALDEHYDE	10.0	9.82	98.23	90-110
15 M,P-TOLUALDEHYDE	20.0	19.5	97.74	90-110
16 HEXANAL	10.0	9.86	98.65	90-110
17 2,5-Dimethylbenzal	10.0	10.1	100.95	90-110

Data File: /chem/hp1cf.i/09Nov2010.b/f1109027.d

Date : 09-NOV-2010 17:40

Client ID: CCV

Sample Info: f1838-41-10;CCV

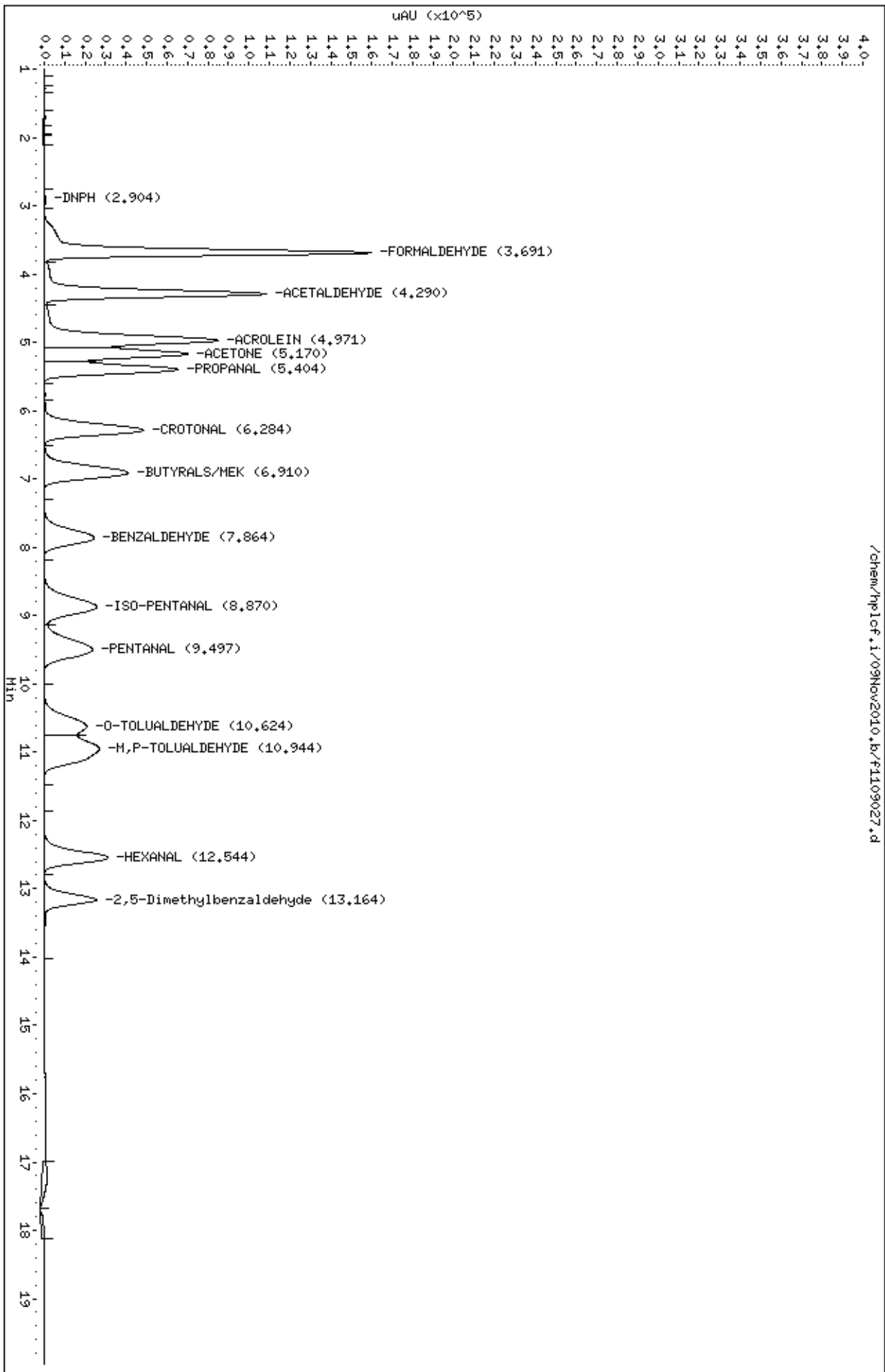
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109028.d
Lab Smp Id: CCB Client Smp ID: LAB BLANK
Inj Date : 09-NOV-2010 18:01
Operator : CRL Inst ID: hplcf.i
Smp Info : ;CCB;LAB BLANK
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	5.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	CONCENTRATIONS	
							ON-COLUMN	FINAL
							(ng)	(ug)
1 DNPH						Compound Not Detected.		
2 FORMALDEHYDE						Compound Not Detected.		
3 ACETALDEHYDE						Compound Not Detected.		
4 ACROLEIN						Compound Not Detected.		
5 ACETONE						Compound Not Detected.		
6 PROPANAL						Compound Not Detected.		
7 CROTONAL						Compound Not Detected.		
9 BUTYRALS/MEK						Compound Not Detected.		
10 BENZALDEHYDE						Compound Not Detected.		
11 ISO-PENTANAL						Compound Not Detected.		
12 PENTANAL						Compound Not Detected.		
14 O-TOLUALDEHYDE						Compound Not Detected.		
15 M,P-TOLUALDEHYDE						Compound Not Detected.		
16 HEXANAL						Compound Not Detected.		

Compounds	RT	CONCENTRATIONS			RESPONSE	ON-COLUMN		FINAL
		EXP	RT	DLT		(	ng)	(
=====	==	=====	=====		=====	=====		=====
17 2,5-Dimethylbenzaldehyde					Compound Not Detected.			

Data File: /chem/hplcf.i/09Nov2010.b/f1109028.d

Page 1

Date : 09-NOV-2010 18:01

Client ID: LAB BLANK

Instrument: hplcf.i

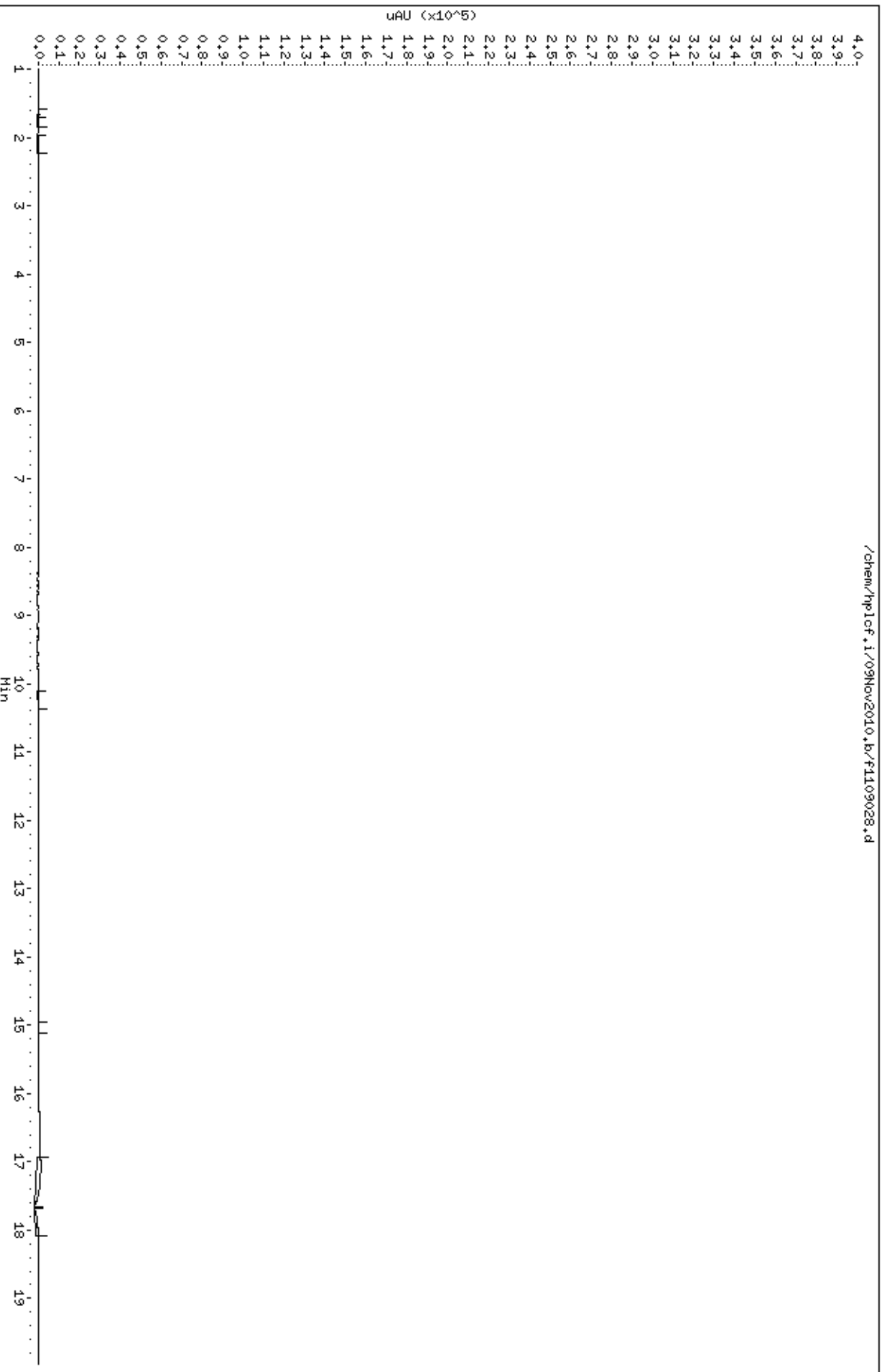
Sample Info: JCCB;LAB BLANK

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

TARGET COMPOUNDS

Client Name:	Client SDG: 09Nov2010
Lab Smp Id: 1838-41-10	Client Smp ID: CCV
Sample Location:	Sample Point:
Sample Date:	Date Received:
Sample Matrix: AIR	Quant Type: ESTD
Analysis Type: OTHER	Level: MED
Data Type: LC DATA	Operator: CRL
Misc Info:	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/KG) ug	Q
119-26-6-----	DNPH	0.00	
50-00-0-----	FORMALDEHYDE	10.5	
75-07-0-----	ACETALDEHYDE	10.0	
107-02-8-----	ACROLEIN	10.5	
67-64-1-----	ACETONE	9.65	
123-38-6-----	PROPANAL	9.58	
123-73-9-----	CROTONAL	9.84	
9999-9999-007--	BUTYRALS/MEK	9.82	
100-52-7-----	BENZALDEHYDE	9.77	
590-86-3-----	ISO-PENTANAL	9.96	
110-62-3-----	PENTANAL	9.82	
529-20-4-----	O-TOLUALDEHYDE	9.97	
620-23-51-----	M, P-TOLUALDEHYDE	19.3	
66-25-1-----	HEXANAL	9.89	
5779-94-2-----	2,5-Dimethylbenzaldehyde	9.96	
=====	=====	=====	=====

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109039.d
Lab Smp Id: 1838-41-10 Client Smp ID: CCV
Inj Date : 09-NOV-2010 21:51
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1838-41-10;CCV
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1 QC Sample: METHSPIKE
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	5.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds					CONCENTRATIONS	
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.663	3.731	-0.068	1148383	10.4900	10.5
3 ACETALDEHYDE	4.257	4.344	-0.087	853629	10.0240	10.0
4 ACRROLEIN	4.923	5.051	-0.128	854203	10.5286	10.5
5 ACETONE	5.123	5.250	-0.127	601489	9.65097	9.65
6 PROPANAL	5.350	5.497	-0.147	600267	9.57607	9.58
7 CROTONAL	6.203	6.417	-0.214	560316	9.83976	9.84
9 BUTYRALS/MEK	6.817	7.064	-0.247	531809	9.82407	9.82
10 BENZALDEHYDE	7.723	8.064	-0.341	373851	9.77164	9.77
11 ISO-PENTANAL	8.710	9.104	-0.394	452769	9.96059	9.96
12 PENTANAL	9.303	9.764	-0.461	438841	9.81743	9.82
14 O-TOLUALDEHYDE	10.430	10.877	-0.447	363534	9.96910	9.97
15 M,P-TOLUALDEHYDE	10.763	11.184	-0.421	591558	19.3107	19.3
16 HEXANAL	12.403	12.717	-0.314	398789	9.88716	9.89
17 2,5-Dimethylbenzaldehyde	13.037	13.324	-0.287	293716	9.95819	9.96

Air Toxics Ltd.

RECOVERY REPORT

Client Name: Client SDG: 09Nov2010
Sample Matrix: LIQUID Fraction: OTHER
Lab Smp Id: 1838-41-10 Client Smp ID: CCV
Level: MED Operator: CRL
Data Type: LC DATA SampleType: METHSPIKE
SpikeList File: 10ug.spk Quant Type: ESTD
Sublist File: all.sub
Method File: /chem/hplcf.i/09Nov2010.b/F10D0914.m
Misc Info:

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	10.0	10.5	104.90	90-110
3 ACETALDEHYDE	10.0	10.0	100.24	90-110
4 ACROLEIN	10.0	10.5	105.29	90-110
5 ACETONE	10.0	9.65	96.51	90-110
6 PROPANAL	10.0	9.58	95.76	90-110
7 CROTONAL	10.0	9.84	98.40	90-110
9 BUTYRALS/MEK	10.0	9.82	98.24	90-110
10 BENZALDEHYDE	10.0	9.77	97.72	90-110
11 ISO-PENTANAL	10.0	9.96	99.61	90-110
12 PENTANAL	10.0	9.82	98.17	90-110
14 O-TOLUALDEHYDE	10.0	9.97	99.69	90-110
15 M,P-TOLUALDEHYDE	20.0	19.3	96.55	90-110
16 HEXANAL	10.0	9.89	98.87	90-110
17 2,5-Dimethylbenzal	10.0	9.96	99.58	90-110

Data File: /chem/hplcf,i/09Nov2010,b/f1109039.d

Date : 09-NOV-2010 21:51

Client ID: CCV

Sample Info: f1838-41-10;CCV

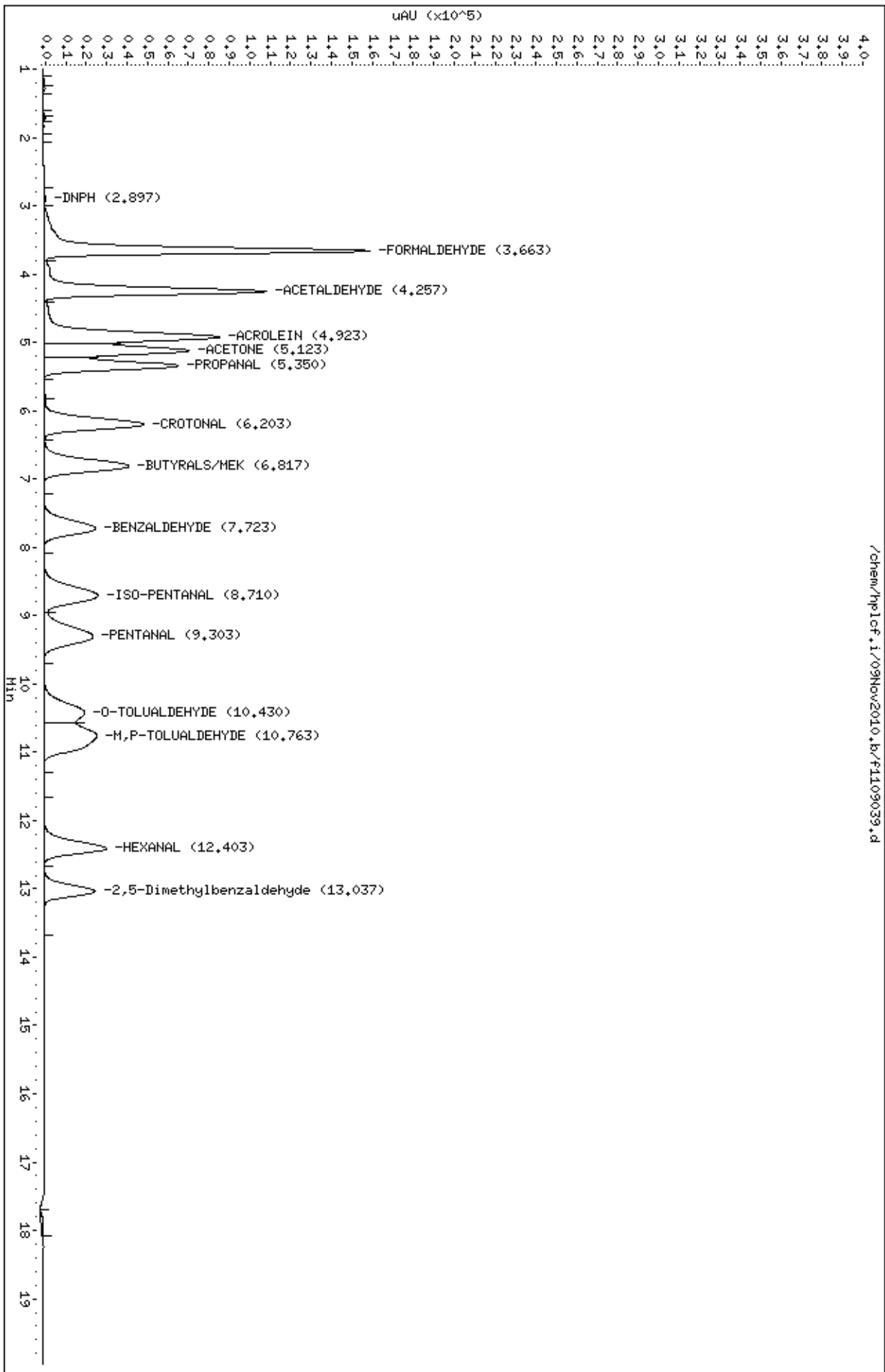
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hplcf,i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109040.d
Lab Smp Id: CCB Client Smp ID: LAB BLANK
Inj Date : 09-NOV-2010 22:12
Operator : CRL Inst ID: hplcf.i
Smp Info : ;CCB;LAB BLANK
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: all.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	5.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	CONCENTRATIONS	
							ON-COLUMN	FINAL
							(ng)	(ug)
1 DNPH						Compound Not Detected.		
2 FORMALDEHYDE						Compound Not Detected.		
3 ACETALDEHYDE						Compound Not Detected.		
4 ACROLEIN						Compound Not Detected.		
5 ACETONE						Compound Not Detected.		
6 PROPANAL						Compound Not Detected.		
7 CROTONAL						Compound Not Detected.		
9 BUTYRALS/MEK						Compound Not Detected.		
10 BENZALDEHYDE						Compound Not Detected.		
11 ISO-PENTANAL						Compound Not Detected.		
12 PENTANAL						Compound Not Detected.		
14 O-TOLUALDEHYDE						Compound Not Detected.		
15 M,P-TOLUALDEHYDE						Compound Not Detected.		
16 HEXANAL						Compound Not Detected.		

Compounds	RT	EXP	RT	DLT	RT	RESPONSE	CONCENTRATIONS	
							ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====			=====	=====	
17 2,5-Dimethylbenzaldehyde		Compound Not Detected.						

Data File: /chem/hp1cf.i/09Nov2010.b/f1109040.d

Page 1

Date : 09-NOV-2010 22:12

Client ID: LAB BLANK

Instrument: hp1cf.i

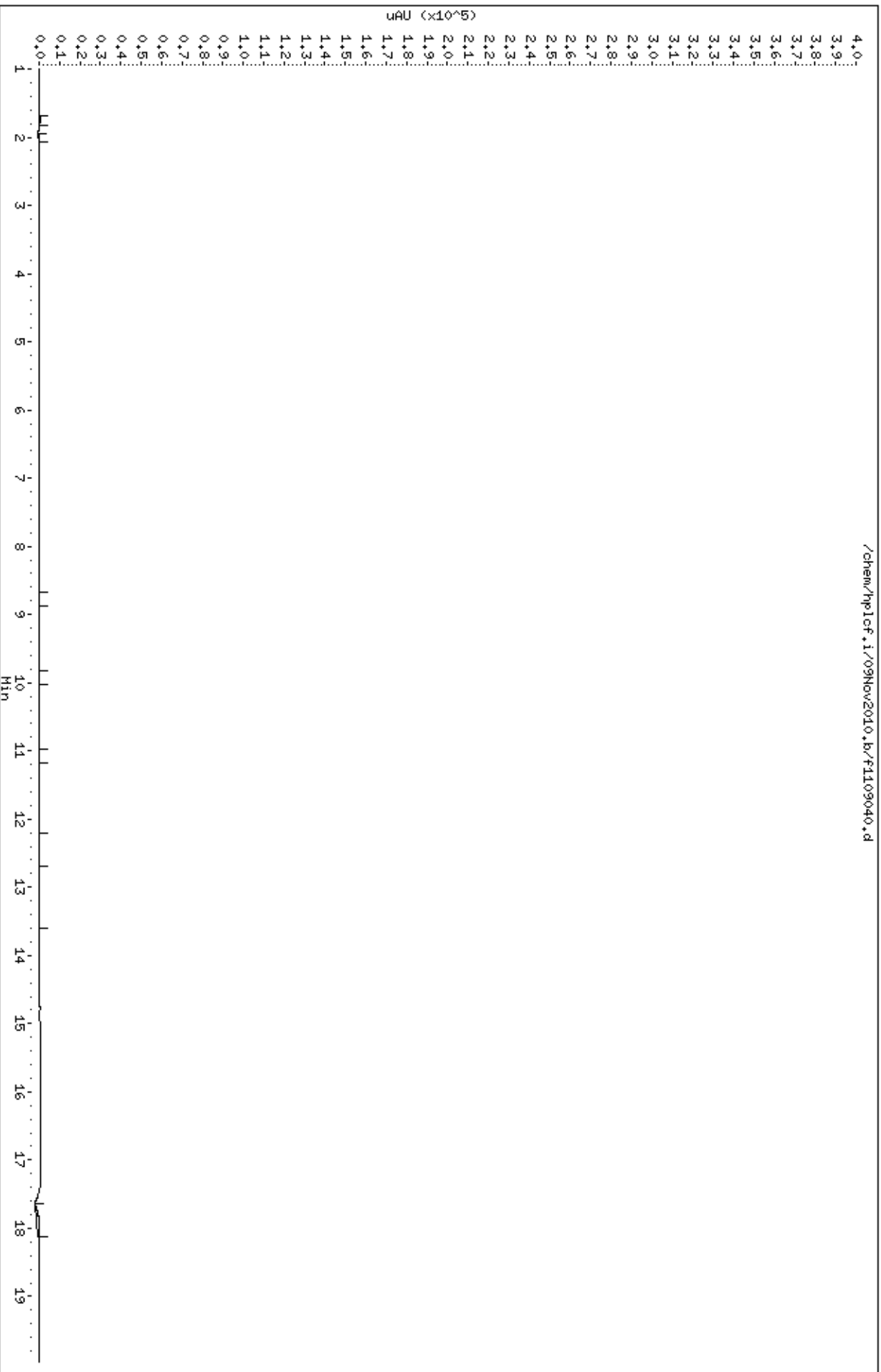
Sample Info: JCCB;LAB BLANK

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

TARGET COMPOUNDS

Client Name:	Client SDG: 09Nov2010
Lab Smp Id: 1838-41-10	Client Smp ID: CCV
Sample Location:	Sample Point:
Sample Date:	Date Received:
Sample Matrix: AIR	Quant Type: ESTD
Analysis Type: OTHER	Level: MED
Data Type: LC DATA	Operator: CRL
Misc Info:	

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/KG) ug	Q
119-26-6-----	DNPH	0.00	
50-00-0-----	FORMALDEHYDE	10.5	
75-07-0-----	ACETALDEHYDE	10.1	
107-02-8-----	ACROLEIN	10.3	
67-64-1-----	ACETONE	9.92	
123-38-6-----	PROPANAL	10.0	
123-73-9-----	CROTONAL	9.99	
9999-9999-007--	BUTYRALS/MEK	10.0	
100-52-7-----	BENZALDEHYDE	10.1	
590-86-3-----	ISO-PENTANAL	10.1	
110-62-3-----	PENTANAL	10.1	
529-20-4-----	O-TOLUALDEHYDE	9.92	
620-23-51-----	M, P-TOLUALDEHYDE	20.0	
66-25-1-----	HEXANAL	9.88	
5779-94-2-----	2,5-Dimethylbenzaldehyde	9.68	
=====	=====	=====	=====

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
 Data file : /chem/hplcf.i/09Nov2010.b/f1109046.d
 Lab Smp Id: 1838-41-10 Client Smp ID: CCV
 Inj Date : 10-NOV-2010 00:17
 Operator : CRL Inst ID: hplcf.i
 Smp Info : ;1838-41-10;CCV
 Misc Info :
 Comment :
 Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
 Meth Date : 09-Nov-2010 11:51 cleaf Quant Type: ESTD
 Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
 Als bottle: 1 QC Sample: METHSPIKE
 Dil Factor: 1.00000
 Integrator: Falcon Compound Sublist: all.sub
 Target Version: 3.50
 Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	5.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds	RT	EXP RT	DLT RT	RESPONSE	CONCENTRATIONS	
					ON-COLUMN (ng)	FINAL (ug)
2 FORMALDEHYDE	3.723	3.731	-0.008	1154199	10.5431	10.5
3 ACETALDEHYDE	4.349	4.344	0.005	859612	10.0942	10.1
4 ACRROLEIN	5.056	5.051	0.005	833913	10.2785	10.3
5 ACETONE	5.249	5.250	-0.001	618372	9.92186	9.92
6 PROPANAL	5.489	5.497	-0.008	627088	10.0039	10.0
7 CROTONAL	6.409	6.417	-0.008	568768	9.98818	9.99
9 BUTYRALS/MEK	7.056	7.064	-0.008	543749	10.0446	10.0
10 BENZALDEHYDE	8.049	8.064	-0.015	385842	10.0850	10.1
11 ISO-PENTANAL	9.083	9.104	-0.021	459804	10.1153	10.1
12 PENTANAL	9.736	9.764	-0.028	450329	10.0744	10.1
14 O-TOLUALDEHYDE	10.856	10.877	-0.021	361935	9.92524	9.92
15 M,P-TOLUALDEHYDE	11.169	11.184	-0.015	611427	19.9593	20.0
16 HEXANAL	12.709	12.717	-0.008	398628	9.88319	9.88
17 2,5-Dimethylbenzaldehyde	13.316	13.324	-0.008	285677	9.68564	9.68

Air Toxics Ltd.

RECOVERY REPORT

Client Name: Client SDG: 09Nov2010
Sample Matrix: LIQUID Fraction: OTHER
Lab Smp Id: 1838-41-10 Client Smp ID: CCV
Level: MED Operator: CRL
Data Type: LC DATA SampleType: METHSPIKE
SpikeList File: 10ug.spk Quant Type: ESTD
Sublist File: all.sub
Method File: /chem/hplcf.i/09Nov2010.b/F10D0914.m
Misc Info:

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	10.0	10.5	105.43	90-110
3 ACETALDEHYDE	10.0	10.1	100.94	90-110
4 ACROLEIN	10.0	10.3	102.79	90-110
5 ACETONE	10.0	9.92	99.22	90-110
6 PROPANAL	10.0	10.0	100.04	90-110
7 CROTONAL	10.0	9.99	99.88	90-110
9 BUTYRALS/MEK	10.0	10.0	100.45	90-110
10 BENZALDEHYDE	10.0	10.1	100.85	90-110
11 ISO-PENTANAL	10.0	10.1	101.15	90-110
12 PENTANAL	10.0	10.1	100.74	90-110
14 O-TOLUALDEHYDE	10.0	9.92	99.25	90-110
15 M,P-TOLUALDEHYDE	20.0	20.0	99.80	90-110
16 HEXANAL	10.0	9.88	98.83	90-110
17 2,5-Dimethylbenzal	10.0	9.68	96.86	90-110

Data File: /chem/hp1cf.i/09Nov2010.b/f1109046.d

Date : 10-NOV-2010 00:17

Client ID: CCV

Sample Info: #1838-41-10;CCV

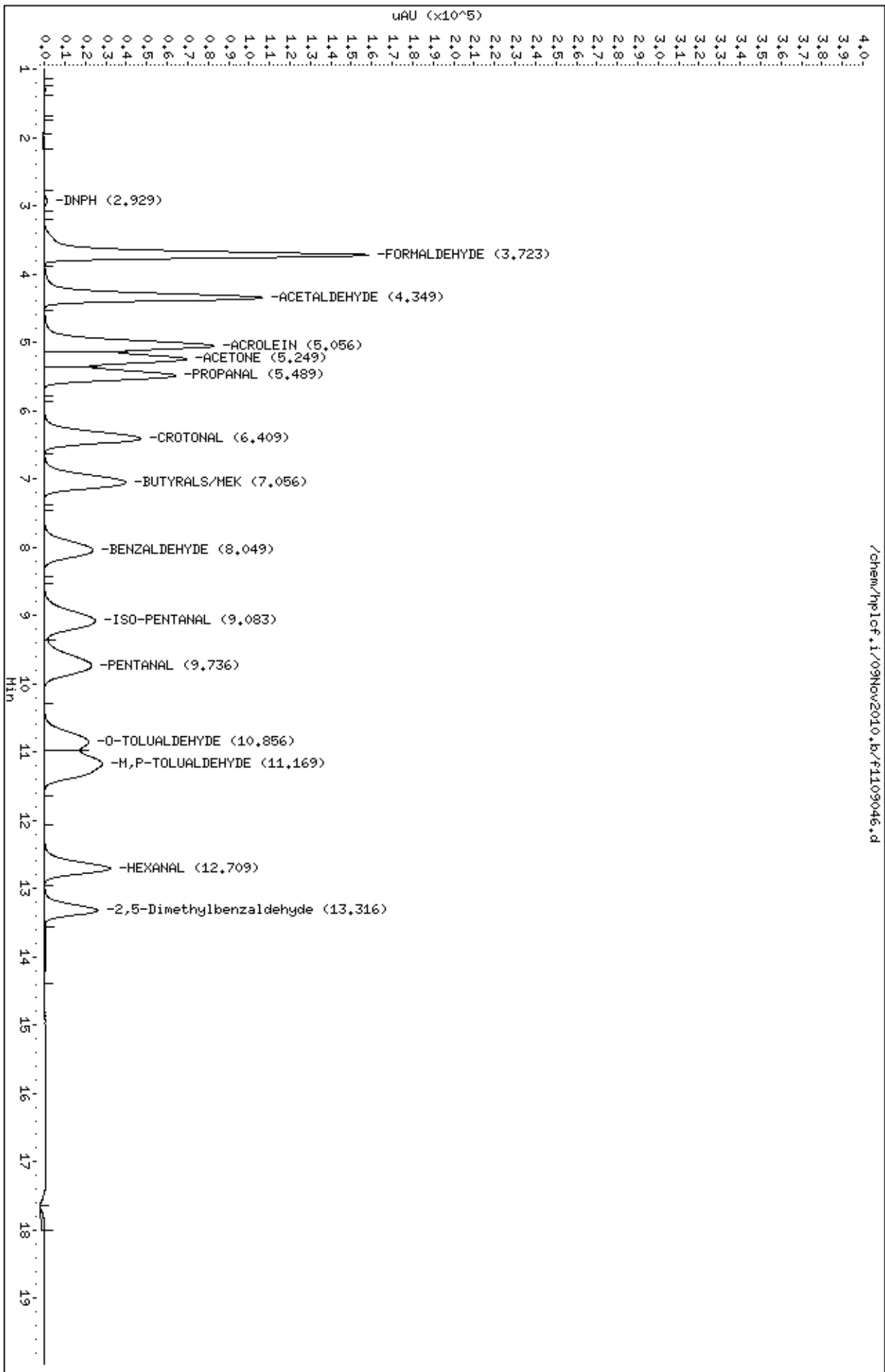
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Client Sample ID: LCS

Lab ID#: 1011114B-27A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1109011
Dil. Factor: 1.00

Date of Collection: NA
Date of Analysis: 11/9/10 12:06 PM
Date of Extraction: 11/9/10

Compound	%Recovery
Formaldehyde	117 Q

Q = Exceeds Quality Control limits.

Container Type: NA - Not Applicable

Air Toxics Ltd.

RECOVERY REPORT

Client Name: Client SDG: 09Nov2010
Sample Matrix: LIQUID Fraction: OTHER
Lab Smp Id: 1011114B Client Smp ID: LCS
Level: MED Operator: CRL
Data Type: LC DATA SampleType: LCS
SpikeList File: FormT0-11A15.spk Quant Type: ESTD
Sublist File: form.sub
Method File: /chem/hplcf.i/09Nov2010.b/F10D0914.m
Misc Info:

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	15.0	17.5	116.77*	85-115

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109011.d
Lab Smp Id: 1011114B Client Smp ID: LCS
Inj Date : 09-NOV-2010 12:06
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1011114B;LCS
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 17-Nov-2010 15:31 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1 QC Sample: LCS
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: form.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	2.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					CONCENTRATIONS	
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.918	2.957	-0.039	32138528		
2 FORMALDEHYDE	3.732	3.731	0.001	4793640	43.7878	17.5(R)

QC Flag Legend

R - Spike/Surrogate failed recovery limits.

Data File: /chem/hp1cf.i/09Nov2010.b/f1109011.d

Page 1

Date : 09-NOV-2010 12:06

Client ID: LCS

Instrument: hp1cf.i

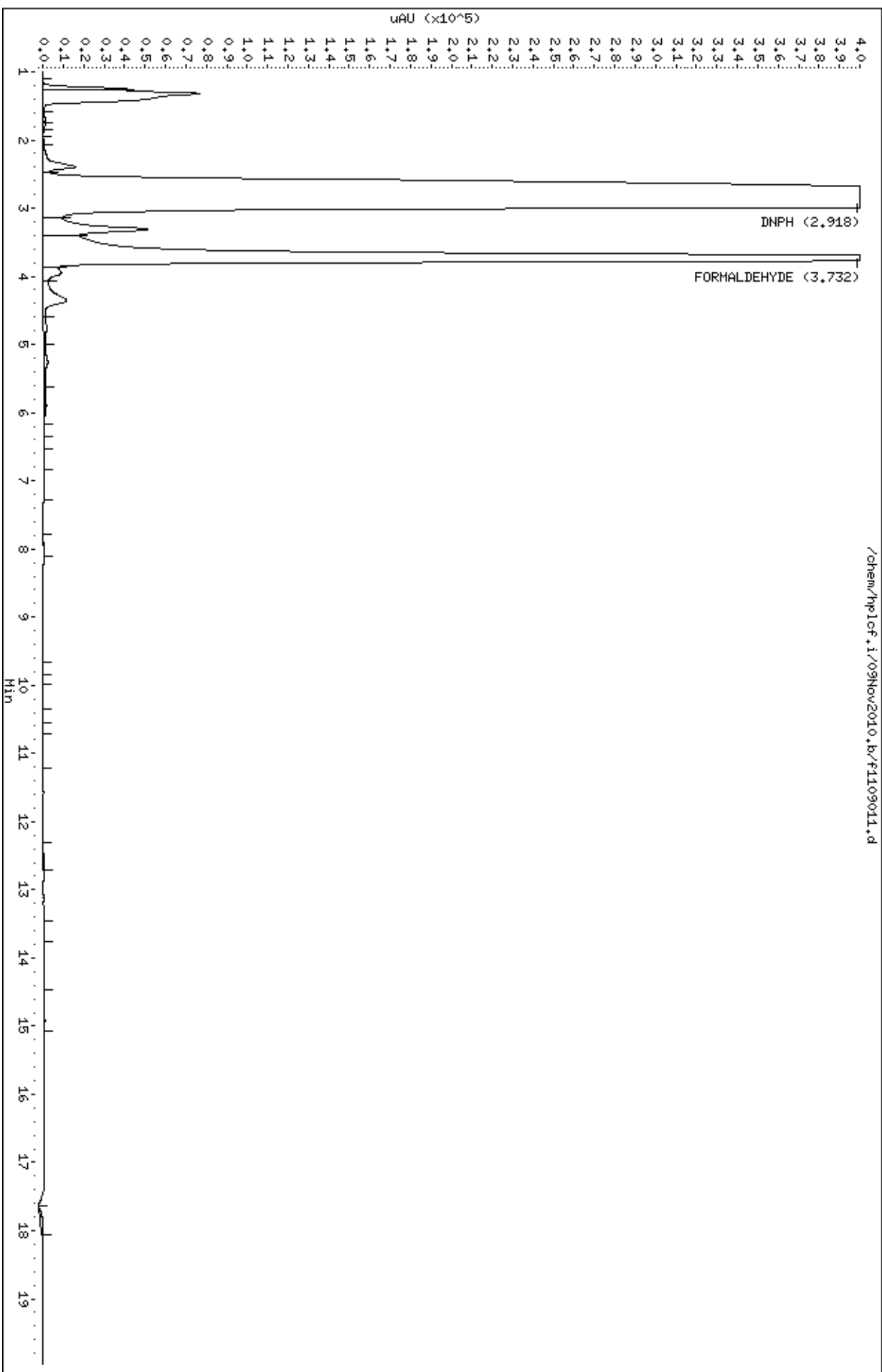
Sample Info: f1011114B:LCS

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Client Sample ID: LCSD

Lab ID#: 1011114B-27AA

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1109012
Dil. Factor: 1.00

Date of Collection: NA
Date of Analysis: 11/9/10 12:27 PM
Date of Extraction: 11/9/10

Compound	%Recovery
Formaldehyde	117 Q

Q = Exceeds Quality Control limits.

Container Type: NA - Not Applicable

Air Toxics Ltd.

RECOVERY REPORT

Client Name: Client SDG: 09Nov2010
Sample Matrix: LIQUID Fraction: OTHER
Lab Smp Id: 1011114B Client Smp ID: LCSD
Level: MED Operator: CRL
Data Type: LC DATA SampleType: LCSD
SpikeList File: FormT0-11A15.spk Quant Type: ESTD
Sublist File: form.sub
Method File: /chem/hplcf.i/09Nov2010.b/F10D0914.m
Misc Info:

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	15.0	17.5	116.67*	85-115

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165
Data file : /chem/hplcf.i/09Nov2010.b/f1109012.d
Lab Smp Id: 1011114B Client Smp ID: LCSD
Inj Date : 09-NOV-2010 12:27
Operator : CRL Inst ID: hplcf.i
Smp Info : ;1011114B;LCSD
Misc Info :
Comment :
Method : /chem/hplcf.i/09Nov2010.b/F10D0914.m
Meth Date : 17-Nov-2010 15:31 cleaf Quant Type: ESTD
Cal Date : 14-SEP-2010 18:30 Cal File: f0914014.d
Als bottle: 1 QC Sample: LCSD
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: form.sub
Target Version: 3.50
Processing Host: eeyore

Concentration Formula: Amt * DF * (Vs/Vo)*Uf * CpndVariable

Name	Value	Description
DF	1.00000	Dilution Factor
Uf	1.00000	ng unit correction factor
Vs	2.00000	Sample Final Volume (mL)
Vo	5.00000	Curve based Volume (mL)

Cpnd Variable Local Compound Variable

Compounds					CONCENTRATIONS	
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.927	2.957	-0.030	32140837		
2 FORMALDEHYDE	3.733	3.731	0.002	4789668	43.7515	17.5(R)

QC Flag Legend

R - Spike/Surrogate failed recovery limits.

Data File: /chem/hplcf.i/09Nov2010.b/f1109012.d

Page 1

Date : 09-NOV-2010 12:27

Client ID: LCSD

Instrument: hplcf.i

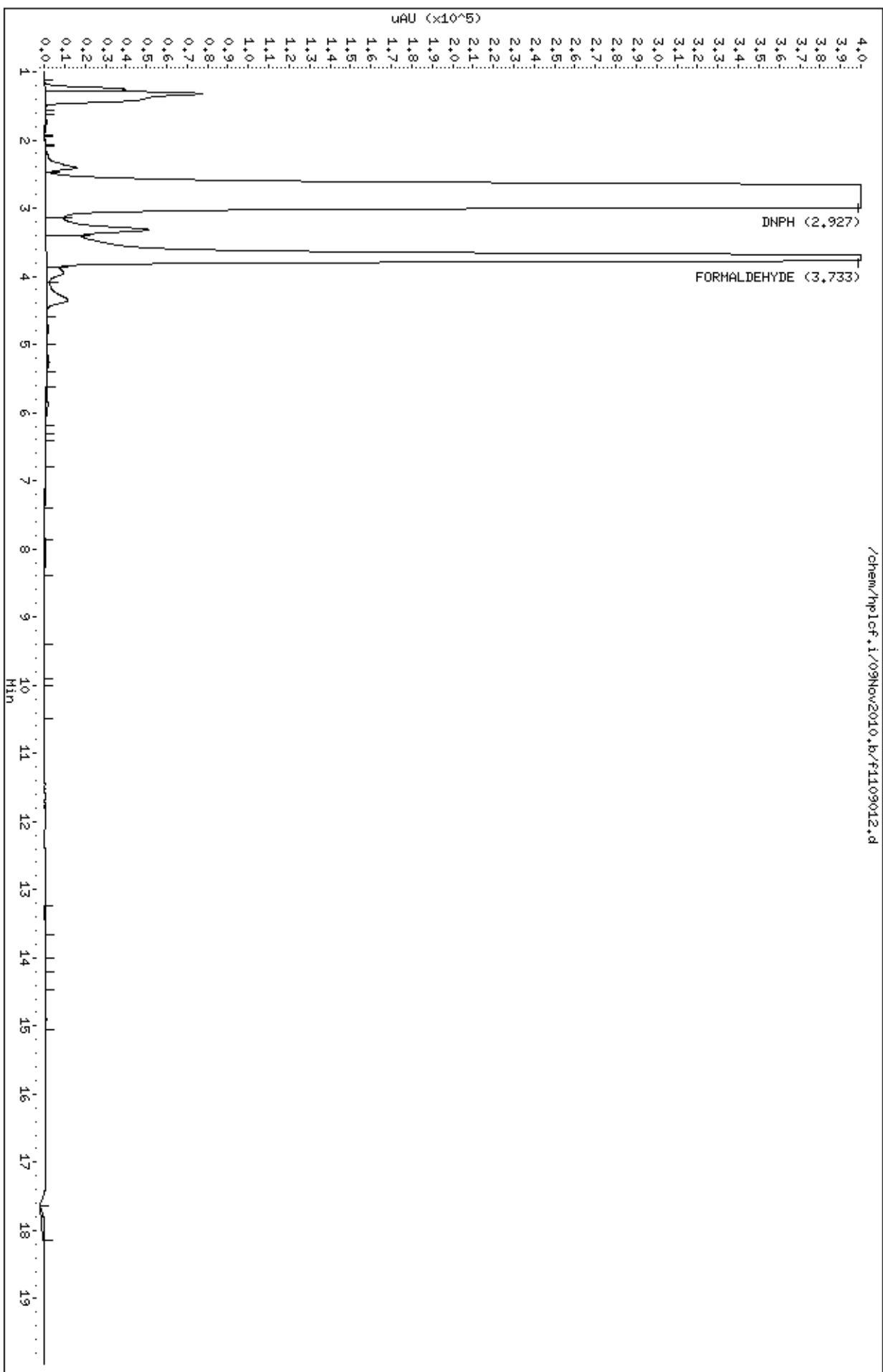
Sample Info: f1011114B;LCSD

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



TO-11 Extraction Logbook

@Air Toxics Ltd.

Log Book #: 2010

ATL Work Order # 101114B

Final Solvent: ACN

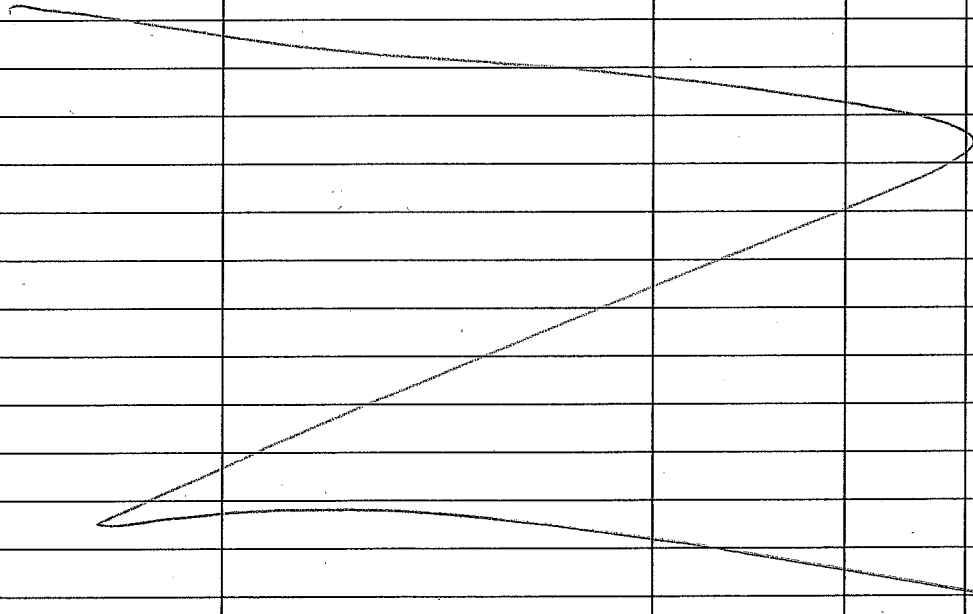
Solvent Lot #: 074016

Date Extracted 11/9/10

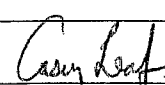
Spike I.D.: 1476493-500ug

Spiked By: CRL

Spike Witness: N/S

ATL Fraction	Clients I.D. #	Cartridge type & size	Final Volume (ml)	Spike Amount (ul)	Comments
101114B-16A	N/A 118635	RAD 165 lot # N/A	2.0	N/A	
-17A	118636				
-18A	118635 118637				
-19A	118636 118638				
-20A	118637 118639				
-21A	118638 118651				
-22A	118639 118652				
-23A	118640 118653				
-24A	118654				
-25A	118655				
101114B-BLK	N/A	lot # 20204			
-UCS		09009		30ul	
					
					Am 11/9/10

Comments:


Signed

11/9/10
Date

Method: Carb 430 / RAD 165Pressure: 122 bar

Use	File #	Sample Name/Client	DF	Loaded by Init.	Date	Time	Reviewed by Init.	Comments
X	1109001	ACN WASH	1.00	CRL	11/9/10	0850 0850	CRL	System shutdown
✓	2	↓				0858 0919		
✓	3	CCV 1838-41-10				0919		
✓	4	CCB				0940		
✓	5	CARB 430 Cert-1				1001		Acetate Form only
✓	6	↓ -2				1022		↓
✓	7	101114A - Lab Blank				1043		
✓	8	↓ -LCS				1103		
✓	9	↓ -LCS				1124		
✓	10	101114B - Lab Blank				1145		
✓	11	↓ -LCS				1206		
✓	12	↓ -LCS				1227		
✓	13	CARB 430 Cert-3				1248		
✓	14	↓ -4				1309		
✓	15	CCV 1838-41-10 mg				1330		
✓	16	CCB				1351		
✓	17	101114A - 05A				1411		
✓	18	↓ -10A				1432		
✓	19	↓ -15A				1453		
✓	20	101114B - 20A				1514		
✓	21	↓ -25A				1535		
✓	22	101114A - 04A				1556		
✓	23	↓ -09A				1617		
✓	24	↓ -14A				1638		
✓	25	101114B - 19A				1659		
✓	26	↓ -24A				1719		
✓	27	CCV 1838-41-10 mg				1740		
✓	28	CCB				1801		

Calculation Check:File ID: 1109014.dCompound: Formal AcetaldehydeInitials: CRL

Sample Amt = $\frac{\text{Area Counts Sample}}{\text{RF}} \times \text{DF} = \left(\frac{17783}{85159} \right) \times (1.00) =$

0.209

Reported Result:

0.209

Pressure: 122 bar

Calculation Check:

Ad 11/10/10

Shipping/ Receiving Documents

Air Toxics Ltd. Sample Receipt Confirmation Cover Page

Thank you for choosing Air Toxics Ltd. We have received your samples and have listed any Sample Receipt Discrepancies below.

In order to expedite analysis and reporting, please review the attached information for
For corrections ca **Ausha Scott at 916-985-1000**

ATL will proceed with the analysis as specified on the Chain of Custody and Sample Receipt Summary page.

Please note : The Sample Receipt Confirmation, including the total workorder charge, is subject to change upon secondary review. Our aim is to provide a confirmation to you in a timely manner. Sample Receipt Discrepancies, if any, may not include discrepancies regarding sample receipt pressure(s). Additionally, the Chain of Custody (COC) will be provided with the final report.

**Environmental
Health &
Engineering, Inc.**

CHAIN OF CUSTODY FORM

DATE: 3 Nov 10

1011114

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

TO: AIR TOXICS

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.)
16A 118635	PASSIVE/AIR	FORMALDEHYDE ANALYSIS 10/20/10-11/2/10	12D 23H 30M
17A 118636			
18A 118637			
19A 118638			
20A 118639			BD
21A 118651		10/19/10-11/2/10	13D 20H 30M
22A 118652			
23A 118653			
24A 118654			
25A 118655			BD

Special instructions:

- ☒ Standard turn around time ☐ Rush by _____ date/time ☐ Other _____
☐ Fax results 781-247-4305
☐ RETURN SAMPLES ☒ Electronic transfer - datacoordinator@ehinc.com
☒ Additional report recipient bb@ehinc.com

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

Relinquished by: [Signature] of Environmental Health & Engineering, Inc. Date: 11/3/10
Received by: Brian Whittaker of (company name) ATL Date: 11/4/10 900
Relinquished by: _____ of (company name) Fedex Date: _____
Received by: _____ of (company name) CUSTODY SEAL INTACT? Y N NONE TEMP 3.5°C Date: _____
Relinquished by: _____ of (company name) Date: _____
Received by: _____ of (company name) Date: _____
Lab Data
Received by: _____ of Environmental Health & Engineering, Inc. Date: _____

Page 2 of 2

SAMPLE RECEIPT SUMMARY

WORKORDER 1011114B

Client Mr. Brian Baker Environmental Health & Engineering, Inc. 117 Fourth Avenue Needham, MA 02494	Phone 800-825-5343 Fax 781-247-4305	Date Promised: 11/18/10 Date Completed: 11/18/10 Date Received: 11/4/10 PO#: 17131 Project#: 17131 Total \$: \$ 950.00 Logged By: MW
Sales Rep: TL		

<u>Fraction</u>	<u>Sample #</u>	<u>Analysis</u>	<u>Collected</u>	<u>Amount\$</u>
16A	118635	Modified TO-11A	11/1/2010	\$90.00
17A	118636	Modified TO-11A	11/1/2010	\$90.00
18A	118637	Modified TO-11A	11/1/2010	\$90.00
19A	118638	Modified TO-11A	11/1/2010	\$90.00
20A	118639	Modified TO-11A	NA	\$90.00
21A	118651	Modified TO-11A	11/1/2010	\$90.00
22A	118652	Modified TO-11A	11/1/2010	\$90.00
23A	118653	Modified TO-11A	11/1/2010	\$90.00
24A	118654	Modified TO-11A	11/1/2010	\$90.00
25A	118655	Modified TO-11A	NA	\$90.00
26A	Lab Blank	Modified TO-11A	NA	\$0.00
27A	LCS	Modified TO-11A	NA	\$0.00
27AA	LCSD	Modified TO-11A	NA	\$0.00
Misc. Charges				
CVP (10) @ \$5.00 each.				\$50.00

Note: Samples received after 3 P.M. PST are considered to be received on the following work day.
 Atlas Project Name/Profile#: CPSC/14482

BILL TO: Accounts Payable
 Environmental Health & Engineering, Inc.
 117 Fourth Avenue
 Needham, MA 02494

Analysis Code: HPLC

TERMS:

Reporting Method: Modified TO-11A Passive-Radiello 165 (Sh)-Form only

180 BLUE RAVINE ROAD, SUITE B FOLSOM, CA - 95630
 (916) 985-1000 . (800) 985-5955 . FAX (916) 985-1020

@ Air Toxics Ltd	Title: Sample Discrepancy Report			Release Date: 03/03/10
	Form #: F1.3	Revision #: 1	Revision Date: 10/7/08	Page #: 1 of 2

Sample Discrepancy Report

Identification

Initiated By: MW Project ID: 14482 PM: As Date: 11/6/2010 Discrepancy Type: ☐ 1. ☒ 2. ☒ 3.

Workorder(s) affected: 1011114 Sample(s) affected: 05A, 10A, 15A, 20A and 25A

1. Sample Receipt Discrepancies

Narration Not Required:

- 1.1. ☐ Sample container (cartridge/tube/VOA vial) was received broken, however sample was intact.
- 1.2. ☐ No brass cap on canister.
- 1.3. ☐ Date of Collection noted on first sample, but no arrow down to indicate all samples.

Notify Lab for further determination:

- 1.4. ☐ Tedlar bag received with minimal volume.

Initials: _____ Date: _____

Narration Required in Lab Narrative and Sample Confirmation:

- 1.5. ☐ COC was not filled out in ink.
- 1.6. ☐ COC improperly relinquished / received.
- 1.7. ☐ Sample tags / can numbers do not match the COC.
- 1.8. ☐ Sample date ☐ error / ☐ missing on COC but noted on sample tag (check one).
- 1.9. ☐ Custody Seal on the outside of the container was ☐ broken / ☐ improperly placed (check one).
- 1.10. ☐ ID-none on the sample Tag/Blank
- 1.11. ☐ Other (describe below).

Describe the Discrepancy: _____

2. Sample Receipt/Screening Discrepancies requiring PM notification

Document on Cover Page of Sample Receipt Confirmation and in Receiving Notes of Lab Narrative

If Section II. is filled out PM must be notified within 24 hrs of initiation

- | | |
|---|---|
| <ol style="list-style-type: none"> 2.1. ☐ COC was not received with samples. 2.2. ☐ Analysis method(s) is ☐ not specified / ☐ incorrectly specified (check one) on the COC. 2.3. ☐ Incorrect sampling media / container for analysis requested. 2.4. ☐ Number of samples on the COC does not match the number of samples that were received. 2.5. ☐ Samples were received expired. 2.6. ☒ Sampling date (time for sulfur) is not documented for ☒ <u>some</u> / ☐ <u>any</u> samples (check one). 2.7. ☐ Sample received with amount of H₂O in the Tedlar Bag. 2.8. ☐ Sample cannot be analyzed. Container was ☐ received broken / ☐ leaking / ☐ flat / ☐ defective. 2.9. ☐ Tedlar bag / canister received emitting a strong odor; Sample ☐ can / ☐ cannot (check one) be analyzed. 2.10. ☐ Tedlar Bag for Sulfur analysis has metal fitting. 2.11. ☐ Environmental Supply Company valves 2.12. ☐ Sorbent samples-sampling volume was not provided | <ol style="list-style-type: none"> 2.13. ☐ Flow controller used – canister samples received at ambient or under pressure. 2.14. ☐ Canister was at ambient pressure at time of pressurization and (check all that apply):
 ☐ Canister failed leak check on two manifolds,
 ☐ Canister valve was open,
 ☐ Brass nut was loose/not present.
 ☐ Sample can be analyzed
 ☐ Cannot be analyzed 2.15. ☐ Canister sample received with a vacuum difference >5.0"Hg between the receipt vac. And the final vac. reported on the COC, indicating loss of vacuum. 2.16. ☐ Canister sample received at >15"Hg (<u>not</u> identified as a Trip/Field Blank). 2.17. ☐ Canister Trip Blank received at low vacuum (< 25"Hg). 2.18. ☐ Sorbent Sample received outside method required temperature of 2°C to 6°C; ☐ ice / ☐ blue ice (check one) was present. A temp. Blank ☐ was / ☐ was not present (check one). 2.19. ☐ Other (describe below) |
|---|---|

Initials

:

Date: _____

Notify Receiving:

☐

Notify PM: ☐

Describe the Discrepancy: No DOCs for samples listed above

3. Lab Discrepancies requiring Team Leader/PM notification

Document in Analytical Notes of Lab Narrative

If Section III. is filled out PM must be notified within 24 hrs of initiation

- | | |
|--|--|
| 3.1. ☐ Tedlar Bag found to be leaking at the time of analysis; sample ☐ can / ☐ cannot (check one) be analyzed. | 3.6. ☐ Sample loss due to instrument malfunction / broken glassware. |
| 3.2. ☐ Tedlar Bag found to be flat/low volume; sample cannot be analyzed. | 3.7. ☐ Low/high surrogate recoveries noted in QC/sample(s) for extractable samples. |
| 3.3. ☐ Sulfur samples received with insufficient time to analyze prior to expiration. | 3.8. ☐ Reporting Limit was raised. |
| 3.4. ☐ Canister found to be leaking at the time of analysis. | 3.9. ☐ Post weight > Pre weight in field/lab Blank for PM10/TSP samples. |
| 3.5. ☐ VOST tube saturated; bag dilution necessary. | 3.10. ☒ Other (describe below). |

Initials

: CRL **Date:** 11/17/2010 **Notify Receiving:** ☐ **Notify PM:** ☐

Team Lead Initials: _____ **Date:** _____

Describe the Discrepancy: 1.) The LCS has recoveries above the acceptance criteria at 116% for Formaldehyde.
The sample was re-run for confirmation. Acceptance criteria for LCS is 85-115%.

How Does this Affect Client:

Project Manager Use Only

Project Manager Notification **Complete**

☒ **Section 2 Complete**

☒ **Section 3**

Action:

- ☒ It is not necessary to notify the client. Narrate the discrepancy in Receiving Notes/Analytical Notes of Lab Narrative.

PM Initials: AS Date: 11/8/2010

- ☐ Client notification required. See attached client contact / email, or comments below:

Client Notification:

PM Initials: _____ Person notified: _____ Date: _____

- ☐ Waiting for Client Reply

Samplers that have "0D" are trip blanks. They will not have DOCs. Please do not discrepancy these.

Comments: _____
3.10 narrate in report

☐ **Notify Lab** Name: _____ Date: _____ **Notify Receiving:** ☐

- ☐ **Additional notifications attached.**

Additional Comments:

Other Records

Compound List

Modified TO-11A Passive-Radiello 165 (Sh)-Form only

CAS Number	Compound	Detection Limit	Type
50-00-0	Formaldehyde	1.0	

@ Air Toxics Ltd	Title: Data Review Checklist		Release Date: 07/28/10	
	Form #: F1.27	Revision #: 2	Revision Date: 07/27/10	Page #: 1 of 2

DATA REVIEW CHECKLIST

Work Order #:

1011143

A ₁	A ₂	W	T	R	Q	
☐	☐	☒	☐	☐	☐	Analysis/Reporting vs. Project Profile/SOP requirements checked (i.e. 100% Dups, J-Flag to MDL, etc)
☐	☐	☒	☐	☐	☐	The final report has the correct reporting list, special units, and header info.
☐	☐	☒	☐	☐	☐	Non-Standard sublist printed/verified, LOQ and LOD verified
☐	☐	☒	☐	☐	☐	Lab Narrative is correct (proper method & description/Receiving & Analytical notes correct)
☐	☐	☒	☐	☐	☐	Sample Discrepancy Report (SDR) is completed
☐	☐	☒	☐	☐	☐	Corrective Action issued - #
☐	☐	☒	☐	☐	☐	Unusual circumstances have been documented in the notes section below
						LUMEN validation report present and initialed
						CIRCLE (YES ☒ NO ☐)
☐	☐	☒	☐	☐	☐	Lab Blank, CCV, LCS and DUP met QC criteria
☐	☐	☒	☐	☐	☐	Hold time is met for all samples
☐	☐	☒	☐	☐	☐	Appropriate data qualifier flags are applied
☐	☐	☒	☐	☐	☐	Manual integrations for samples and QC are properly documented
☐	☐	☒	☐	☐	☐	Samples analyzed within the project or method specific clock
☐	☐	☒	☐	☐	☐	Retention times have been verified
☐	☐	☒	☐	☐	☐	Appropriate ICAL(s) included, %RSD Recalculation
☐	☐	☒	☐	☐	☐	At least one result per sample is verified against the target quant sheets/raw data
☐	☐	☒	☐	☐	☐	Dilution factor correctly calculated (sample load volume, syringe and bag dilutions, can pressurization(s))
☐	☐	☒	☐	☐	☐	Correct amount of sample analyzed (i.e. sample not over-diluted)
☐	☐	☒	☐	☐	☐	Spectra verified - documentation of spectral defense included (Section 5A of eCVP pkg)
☐	☐	☒	☐	☐	☐	TICs resemble reference spectra
☐	☐	☒	☐	☐	☐	TICs between duplicate samples are consistent
☐	☐	☒	☐	☐	☐	Checked samples for trends (i.e. Influent vs. Effluent, Field Dups, Field/Trip Blank, etc.)
☐	☐	☒	☐	☐	☐	Data for multiple analyses of sample(s) has been evaluated for comparability of results
☐	☐	☒	☐	☐	☐	Special units for all samples in the final report are correctly calculated
☐	☐	☒	☐	☐	☐	Manually entered results checked (i.e. TPH/NMOC)
☐	☐	☒	☐	☐	☐	Chain of Custody verified for any special comments (i.e. different compounds/RLs, action levels)
☐	☐	☒	☐	☐	☐	Chain of Custody scanned correctly
☐	☐	☒	☐	☐	☐	Verify sample id's vs. chain of custody
☐	☐	☒	☐	☐	☐	Date MDL(s) performed per instrument(s) 10/17/10
☐	☐	☒	☐	☐	☐	Samples pressurized w/ appropriate gas (N ₂ or He) ☒ Other (i.e. Tedlar bag, cartridge, sorbent)
☐	☐	☒	☐	☐	☐	Final pressure consistent with canister size (6L vs. 1L)
☐	☐	☒	☐	☐	☐	Verify receipt pressures
☐	☐	☒	☐	☐	☐	Verify canister ID #'s
☐	☐	☒	☐	☐	☐	Final invoice amount correct (adjusted for TAT, Penalties, Re-issue Charges etc.)
☐	☐	☒	☐	☐	☐	Final PDF report reviewed for correctness

Notes: (to include: noting samples with QA/QC problems, Blanks with positive hits, narratives, etc.)

A/R: 1) LCS/LSSD = Dup 2) ^{data utilized} LCS out of acceptance criteria @ 116%
 3) FB=20A, 2SA 4) Final volume = 2 or L

T/Q:

A ₁ /A ₂ (Analytical Review/Date)	W/T (Write-up/Tech Review/Date)	R* (Report Review/Date)	Q (QA Review/Date)
A ₁ :	W: <i>Carry Leaf</i> 11/12/10	R:	
A ₂ :	T:		

Note (1): Please check all the appropriate boxes. Indicate "NA" for any statement that does not apply.

Note (2): Report reviewer and write-up reviewer must be separate individuals for DoD & Client Specific projects.

* Report Review is completed for DoD & Client Specific projects only.

Not Applicable