

**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 #: 1011418A

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Comments:

Completed by:

Kara McKiernan

(Signature)

Kara McKiernan / Document Control

(Print Name & Title)

12/7/10

(Date)

WORK ORDER #: 1011418A

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:	11/18/2010	CONTACT:	Ausha Scott
DATE COMPLETED:	12/06/2010		

<u>FRACTION #</u>	<u>NAME</u>	<u>TEST</u>
01A	118683	Modified TO-11A
02A	118684	Modified TO-11A
03A	118685	Modified TO-11A
04A	118686	Modified TO-11A
05A	118687	Modified TO-11A
06A	118699	Modified TO-11A
07A	118700	Modified TO-11A
08A	118701	Modified TO-11A
09A	118702	Modified TO-11A
10A	118703	Modified TO-11A
11A	118715	Modified TO-11A
12A	118716	Modified TO-11A
13A	118717	Modified TO-11A
14A	118718	Modified TO-11A
15A	118719	Modified TO-11A
16A	Lab Blank	Modified TO-11A
17A	LCS	Modified TO-11A

Continued on next page

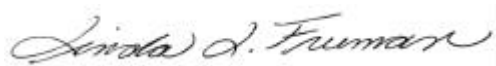
WORK ORDER #: 1011418A

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:	11/18/2010	CONTACT:	Ausha Scott
DATE COMPLETED:	12/06/2010		

<u>FRACTION #</u>	<u>NAME</u>	<u>TEST</u>
17AA	LCSD	Modified TO-11A

CERTIFIED BY:



Laboratory Director

DATE: 12/06/10

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

LABORATORY NARRATIVE
Aldehydes by Radiello 165
Environmental Health & Engineering, Inc.
Workorder# 1011418A

Fifteen Radiello 165 (Aldehydes) samples were received on November 18, 2010. The procedure involves reaction of aldehydes with 2, 4-DNPH to give 2, 4-dinitrohydrazons. The 2, 4-dinitrohydrazons are then extracted with acetonitrile and analyzed using reverse phase High Pressure Liquid Chromatography (HPLC) with an Ultraviolet (UV) Detector. Results are reported in uG/m³.

Receiving Notes

There were no receiving discrepancies.

Analytical Notes

Results were calculated based on 25 deg C without temperature correction. Sampling rates were provided by the manufacturer.

The exposure time for all samples exceeded the method recommended exposure time of 7 days.

The actual exposure time was used to calculate sample concentrations and reporting limits. An exposure time of 20120 minutes was used for the Laboratory Blank.

Definition of Data Qualifying Flags

Eight qualifiers may have been used on the data analysis sheets and indicates as follows:

B - Compound present in laboratory blank greater than reporting limit (background subtraction not performed).

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 reporting limit.

UJ- Non-detected compound associated with low bias in the CCV

N - The identification is based on presumptive evidence.

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
118683	1011418A-01A	11/15/2010	11/18/2010	11/23/2010	8	11/24/2010	1	Good
118684	1011418A-02A	11/15/2010	11/18/2010	11/23/2010	8	11/24/2010	1	Good
118685	1011418A-03A	11/15/2010	11/18/2010	11/23/2010	8	11/24/2010	1	Good
118686	1011418A-04A	11/15/2010	11/18/2010	11/23/2010	8	11/23/2010	0	Good
118687	1011418A-05A	NA	11/18/2010	11/23/2010	NA	11/23/2010	0	Good
118699	1011418A-06A	11/15/2010	11/18/2010	11/23/2010	8	11/24/2010	1	Good
118700	1011418A-07A	11/15/2010	11/18/2010	11/23/2010	8	11/24/2010	1	Good
118701	1011418A-08A	11/15/2010	11/18/2010	11/23/2010	8	11/24/2010	1	Good
118702	1011418A-09A	11/15/2010	11/18/2010	11/23/2010	8	11/23/2010	0	Good
118703	1011418A-10A	NA	11/18/2010	11/23/2010	NA	11/23/2010	0	Good
118715	1011418A-11A	11/15/2010	11/18/2010	11/23/2010	8	11/24/2010	1	Good
118716	1011418A-12A	11/15/2010	11/18/2010	11/23/2010	8	11/24/2010	1	Good
118717	1011418A-13A	11/15/2010	11/18/2010	11/23/2010	8	11/24/2010	1	Good
118718	1011418A-14A	11/15/2010	11/18/2010	11/23/2010	8	11/23/2010	0	Good
118719	1011418A-15A	11/15/2010	11/18/2010	11/23/2010	8	11/23/2010	0	Good
Lab Blank	1011418A-16A	NA	NA	11/23/2010	NA	11/23/2010	0	Good
LCS	1011418A-17A	NA	NA	11/23/2010	NA	11/23/2010	0	Good
LCSD	1011418A-17AA	NA	NA	11/23/2010	NA	11/23/2010	0	Good

Sample Results and Raw Data

Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118683

Lab ID#: 1011418A-01A

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	44	23

Client Sample ID: 118683

Lab ID#: 1011418A-01A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123049
Dil. Factor: 10.0

Date of Collection: 11/15/10
Date of Analysis: 11/24/10 12:22 AM
Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	44	23

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123049.d

Lab Smp Id: 1011418A-01AClient Smp ID: 10

Inj Date : 24-NOV-2010 00:22

Operator : FMInst ID: hplcf.i

Smp Info : ;1011418A-01A;10

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 10.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS					
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.865	2.885	-0.020	2691689		
2 FORMALDEHYDE	3.611	3.598	0.013	1210393	11.0564	44.2

Data File: /chem/hplcf.i/23Nov2010.b/f1123049.d

Page 1

Date : 24-NOV-2010 00:22

Client ID: 10

Instrument: hplcf.i

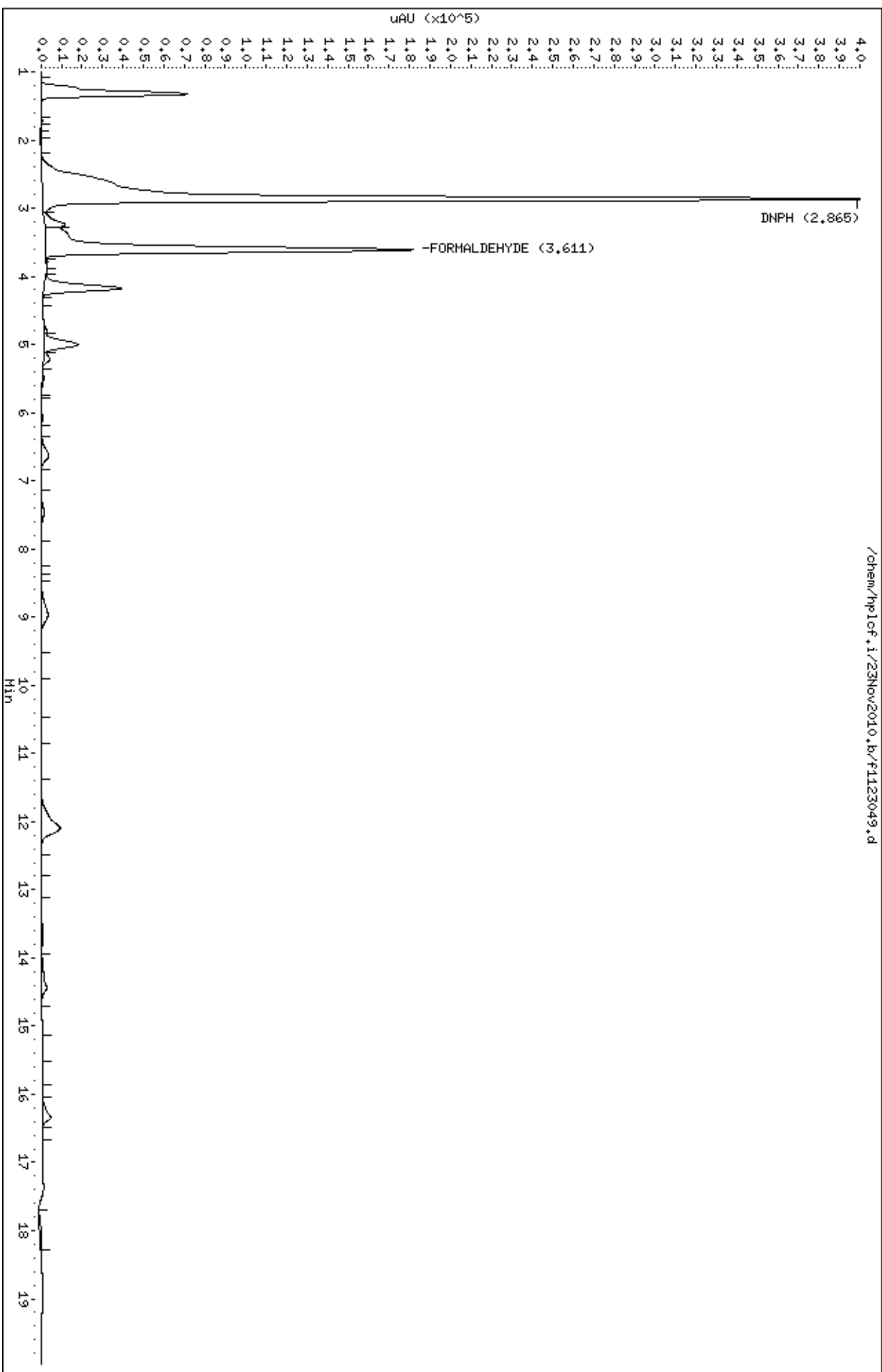
Sample Info: j1011418a-01a;10

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118684

Lab ID#: 1011418A-02A

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	83	43

Client Sample ID: 118684

Lab ID#: 1011418A-02A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123050
Dil. Factor: 10.0

Date of Collection: 11/15/10
Date of Analysis: 11/24/10 12:43 AM
Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	83	43

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123050.d

Lab Smp Id: 1011418A-02AClient Smp ID: 10

Inj Date : 24-NOV-2010 00:43

Operator : FMInst ID: hplcf.i

Smp Info : ;1011418A-02A;10

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 10.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS					
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.788	2.885	-0.097	2436061		
2 FORMALDEHYDE	3.541	3.598	-0.057	2266638	20.7048	82.8

Data File: /chem/hp1cf.i/23Nov2010.b/f1123050.d

Page 1

Date : 24-NOV-2010 00:43

Client ID: 10

Instrument: hp1cf.i

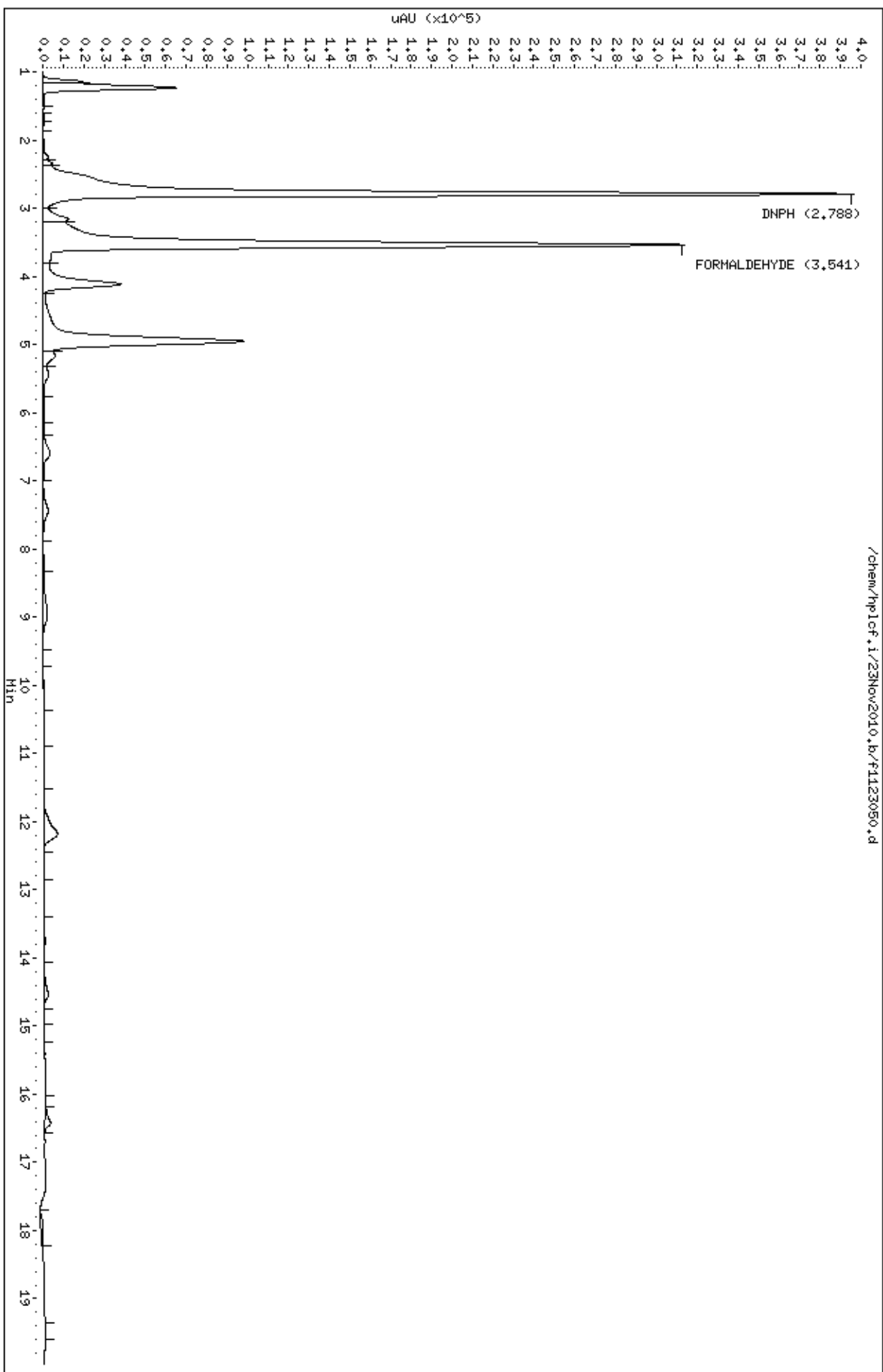
Sample Info: j1011418a-02a;10

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118685

Lab ID#: 1011418A-03A

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	53	27

Client Sample ID: 118685

Lab ID#: 1011418A-03A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123051
Dil. Factor: 10.0

Date of Collection: 11/15/10
Date of Analysis: 11/24/10 01:04 AM
Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	53	27

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123051.d

Lab Smp Id: 1011418A-03AClient Smp ID: 10

Inj Date : 24-NOV-2010 01:04

Operator : FMInst ID: hplcf.i

Smp Info : ;1011418A-03A;10

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 10.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS					
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.853	2.885	-0.032	3015178		
2 FORMALDEHYDE	3.600	3.598	0.002	1445458	13.2036	52.8

Data File: /chem/hp1cf.i/23Nov2010.b/f1123051.d

Page 1

Date : 24-NOV-2010 01:04

Client ID: 10

Instrument: hp1cf.i

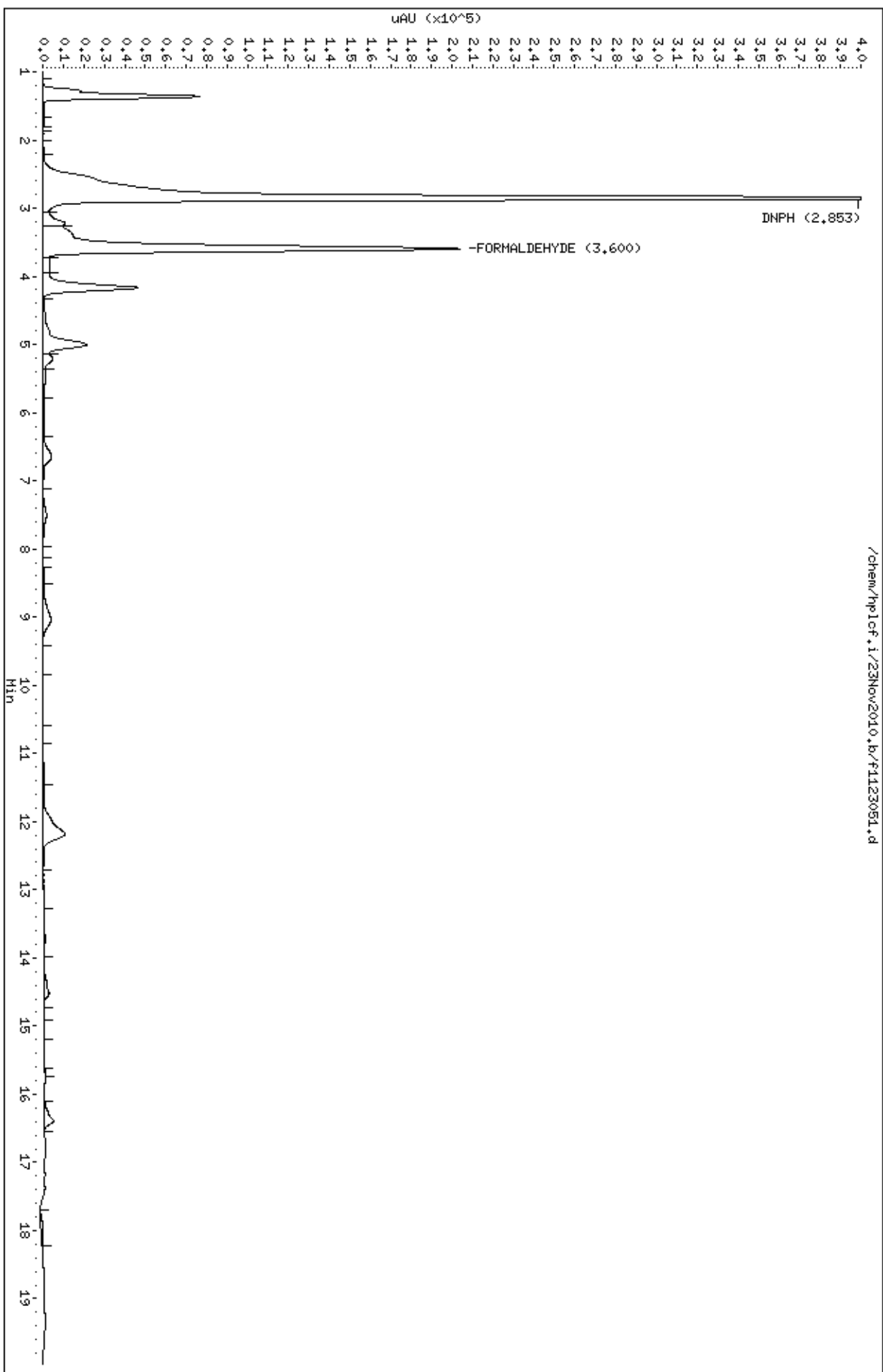
Sample Info: j1011418a-03a;10

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118686

Lab ID#: 1011418A-04A

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	1.0	0.52	2.5	1.3

Client Sample ID: 118686

Lab ID#: 1011418A-04A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123041
Dil. Factor: 1.00

Date of Collection: 11/15/10
Date of Analysis: 11/23/10 09:35 PM
Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	1.0	0.52	2.5	1.3

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123041.d

Lab Smp Id: 1011418A-04A

Inj Date : 23-NOV-2010 21:35

Operator : FMInst ID: hplcf.i

Smp Info : ;1011418A-04A;

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)	
=====	==	=====	=====	=====	=====	=====	
1 DNPH	2.878	2.885	-0.007	15660529			
2 FORMALDEHYDE	3.644	3.598	0.046	686986	6.27533	2.51	

Data File: /chem/hp1cf.i/23Nov2010.b/f1123041.d

Date : 23-NOV-2010 21:35

Client ID:

Sample Info: j1011418a-04a;

Purge Volume: 5.0

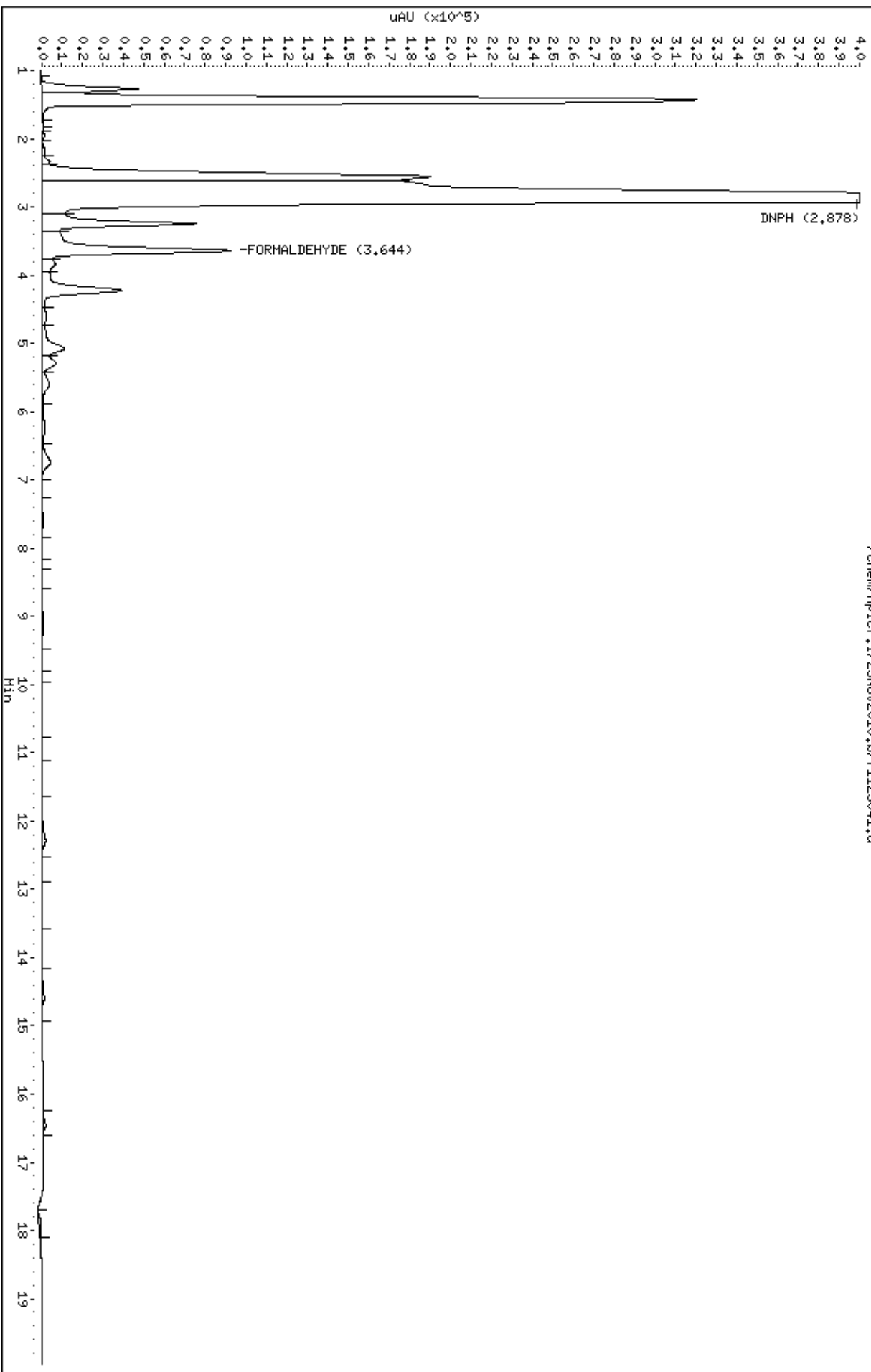
Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: FM

Column diameter: 2.00

/chem/hp1cf.i/23Nov2010.b/f1123041.d



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

Client Sample ID: 118687

Lab ID#: 1011418A-05A

No Detections Were Found.

Client Sample ID: 118687

Lab ID#: 1011418A-05A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123035
Dil. Factor: 1.00

Date of Collection: NA
Date of Analysis: 11/23/10 07:30 PM
Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	1.0	0.52	Not Detected	Not Detected

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123035.d

Lab Smp Id: 1011418A-05A

Inj Date : 23-NOV-2010 19:30

Operator : FMInst ID: hplcf.i

Smp Info : ;1011418A-05A;

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS					
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.885	2.885	0.000	29596801		
2 FORMALDEHYDE				Compound Not Detected.		

Data File: /chem/hp1cf.i/23Nov2010.b/f1123035.d

Page 1

Date : 23-NOV-2010 19:30

Client ID:

Instrument: hp1cf.i

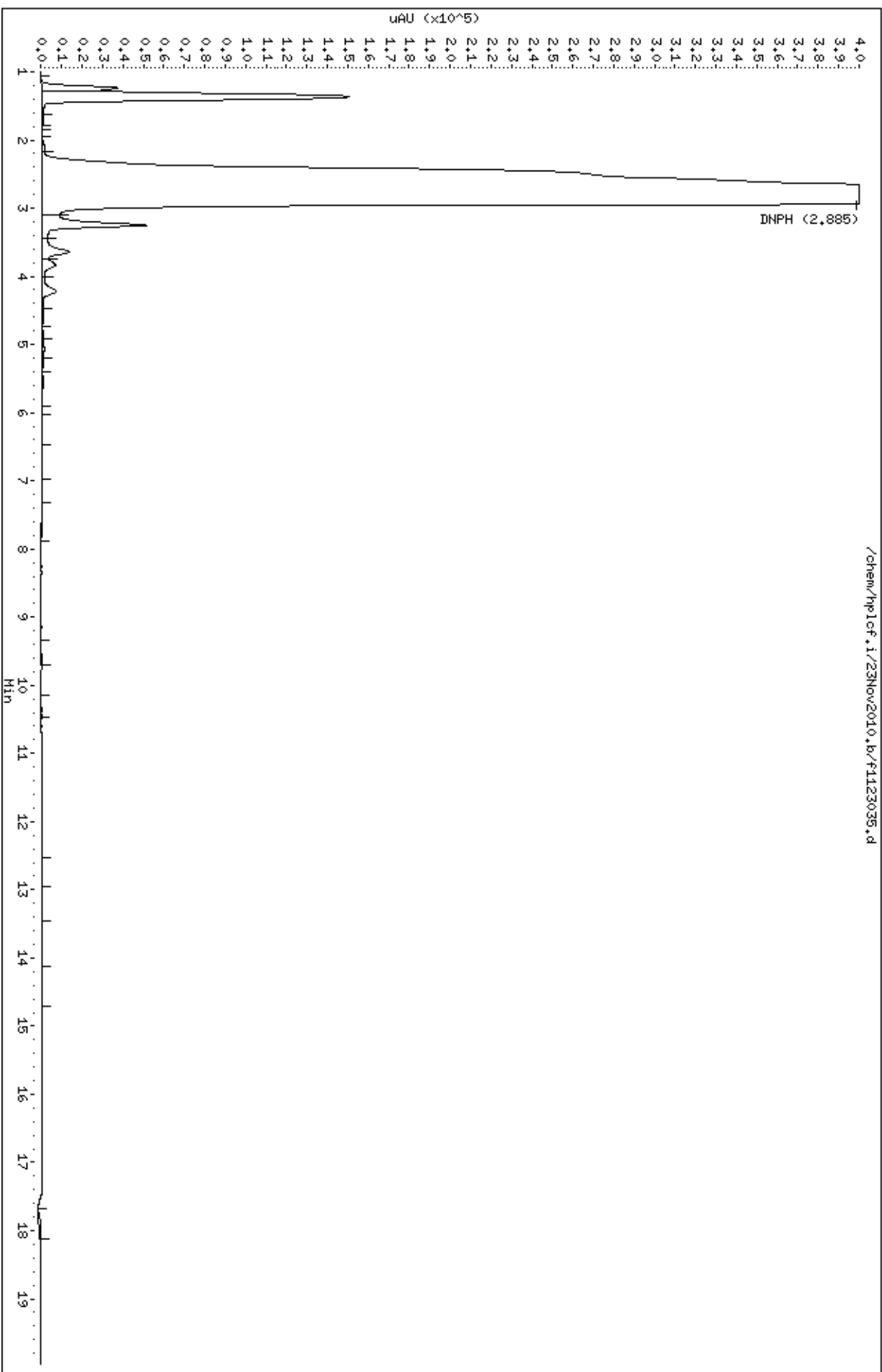
Sample Info: j1011418a-05a;

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118699

Lab ID#: 1011418A-06A

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	70	36

Client Sample ID: 118699

Lab ID#: 1011418A-06A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123052
Dil. Factor: 10.0

Date of Collection: 11/15/10
Date of Analysis: 11/24/10 01:25 AM
Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	70	36

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123052.d

Lab Smp Id: 1011418A-06AClient Smp ID: 10

Inj Date : 24-NOV-2010 01:25

Operator : FMInst ID: hplcf.i

Smp Info : ;1011418A-06A;10

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 10.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS					
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.881	2.885	-0.004	2773491		
2 FORMALDEHYDE	3.641	3.598	0.043	1914285	17.4862	69.9

Data File: /chem/hp1cf.i/23Nov2010.b/f1123052.d

Page 1

Date : 24-NOV-2010 01:25

Client ID: 10

Instrument: hp1cf.i

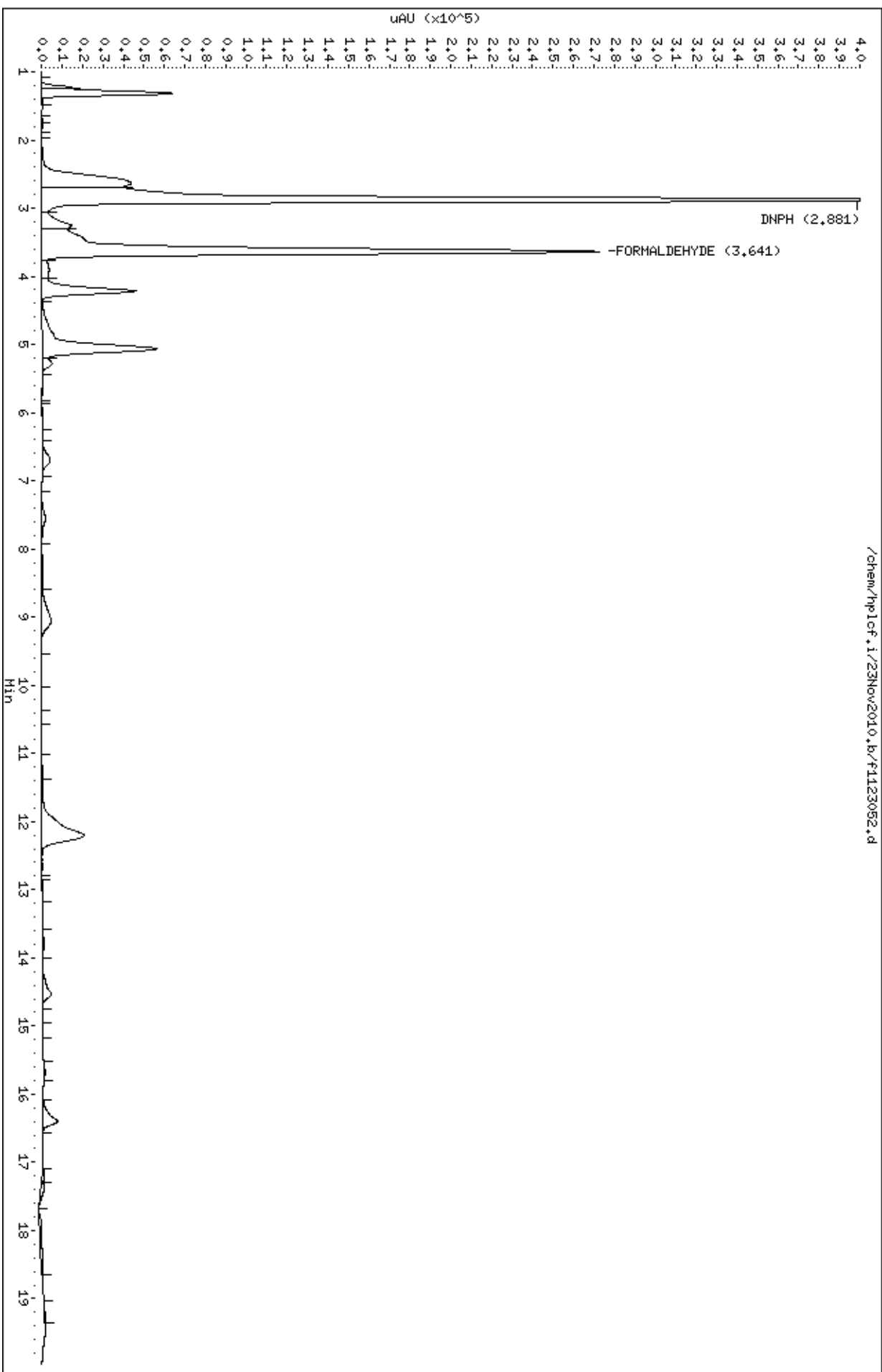
Sample Info: j1011418a-06a;10

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118700

Lab ID#: 1011418A-07A

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	63	33

Client Sample ID: 118700

Lab ID#: 1011418A-07A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123053
Dil. Factor: 10.0

Date of Collection: 11/15/10
Date of Analysis: 11/24/10 01:46 AM
Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	63	33

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123053.d

Lab Smp Id: 1011418A-07AClient Smp ID: 10

Inj Date : 24-NOV-2010 01:46

Operator : FMInst ID: hplcf.i

Smp Info : ;1011418A-07A;10

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 10.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS					
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.841	2.885	-0.044	2158646		
2 FORMALDEHYDE	3.602	3.598	0.004	1734640	15.8452	63.4

Data File: /chem/hp1cf.i/23Nov2010.b/f1123053.d

Page 1

Date : 24-NOV-2010 01:46

Client ID: 10

Instrument: hp1cf.i

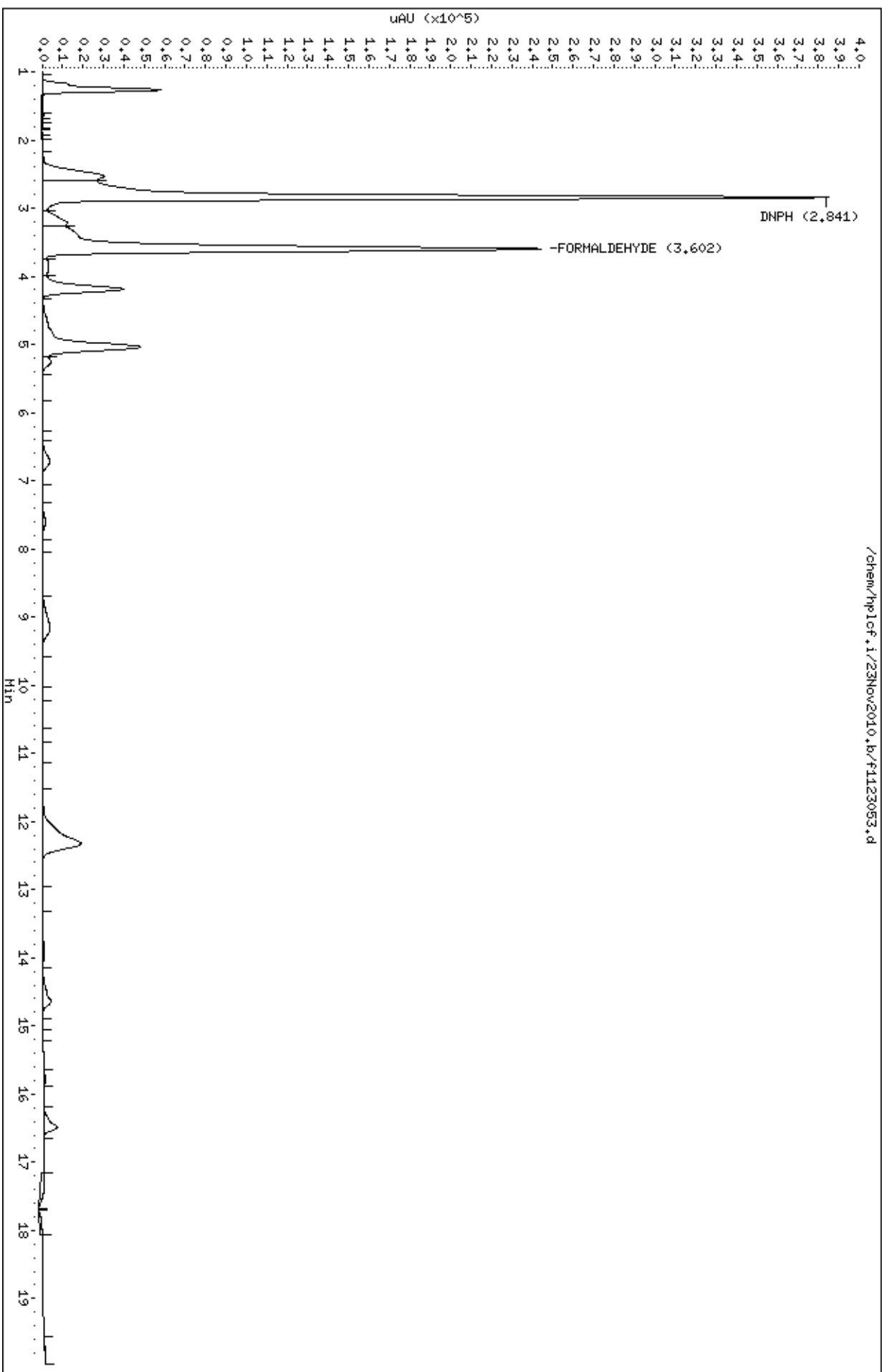
Sample Info: j1011418a-07a;10

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118701

Lab ID#: 1011418A-08A

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	76	39

Client Sample ID: 118701

Lab ID#: 1011418A-08A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123056
Dil. Factor: 10.0

Date of Collection: 11/15/10
Date of Analysis: 11/24/10 02:49 AM
Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	76	39

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123056.d

Lab Smp Id: 1011418A-08AClient Smp ID: 10

Inj Date : 24-NOV-2010 02:49

Operator : FMInst ID: hplcf.i

Smp Info : ;1011418A-08A;10

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 10.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS					
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.847	2.885	-0.038	1548141		
2 FORMALDEHYDE	3.600	3.598	0.002	2083454	19.0314	76.1

Data File: /chem/hp1cf.i/23Nov2010.b/f1123056.d

Page 1

Date : 24-NOV-2010 02:49

Client ID: 10

Instrument: hp1cf.i

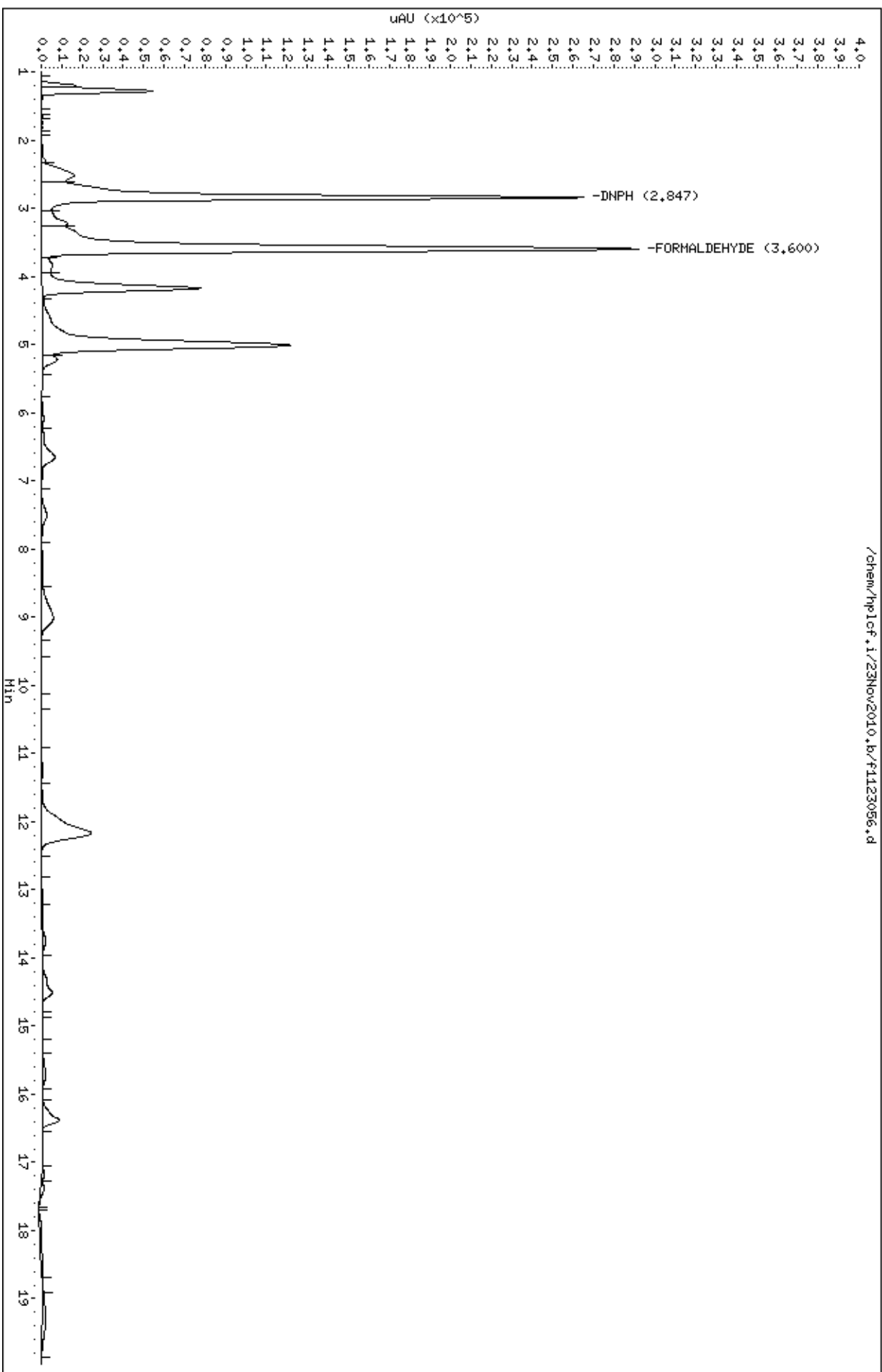
Sample Info: j1011418a-08a;10

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118702

Lab ID#: 1011418A-09A

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	1.0	0.52	2.0	1.0

Client Sample ID: 118702

Lab ID#: 1011418A-09A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123044

Date of Collection: 11/15/10

Dil. Factor: 1.00

Date of Analysis: 11/23/10 10:38 PM

Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	1.0	0.52	2.0	1.0

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123044.d

Lab Smp Id: 1011418A-09A

Inj Date : 23-NOV-2010 22:38

Operator : FM

Smp Info : ;1011418A-09A;

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 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	2.897	2.885	0.012	23632185		
2 FORMALDEHYDE	3.657	3.598	0.059	549723	5.02148	2.01

Data File: /chem/hp1cf.i/23Nov2010.b/f1123044.d

Page 1

Date : 23-NOV-2010 22:38

Client ID:

Instrument: hp1cf.i

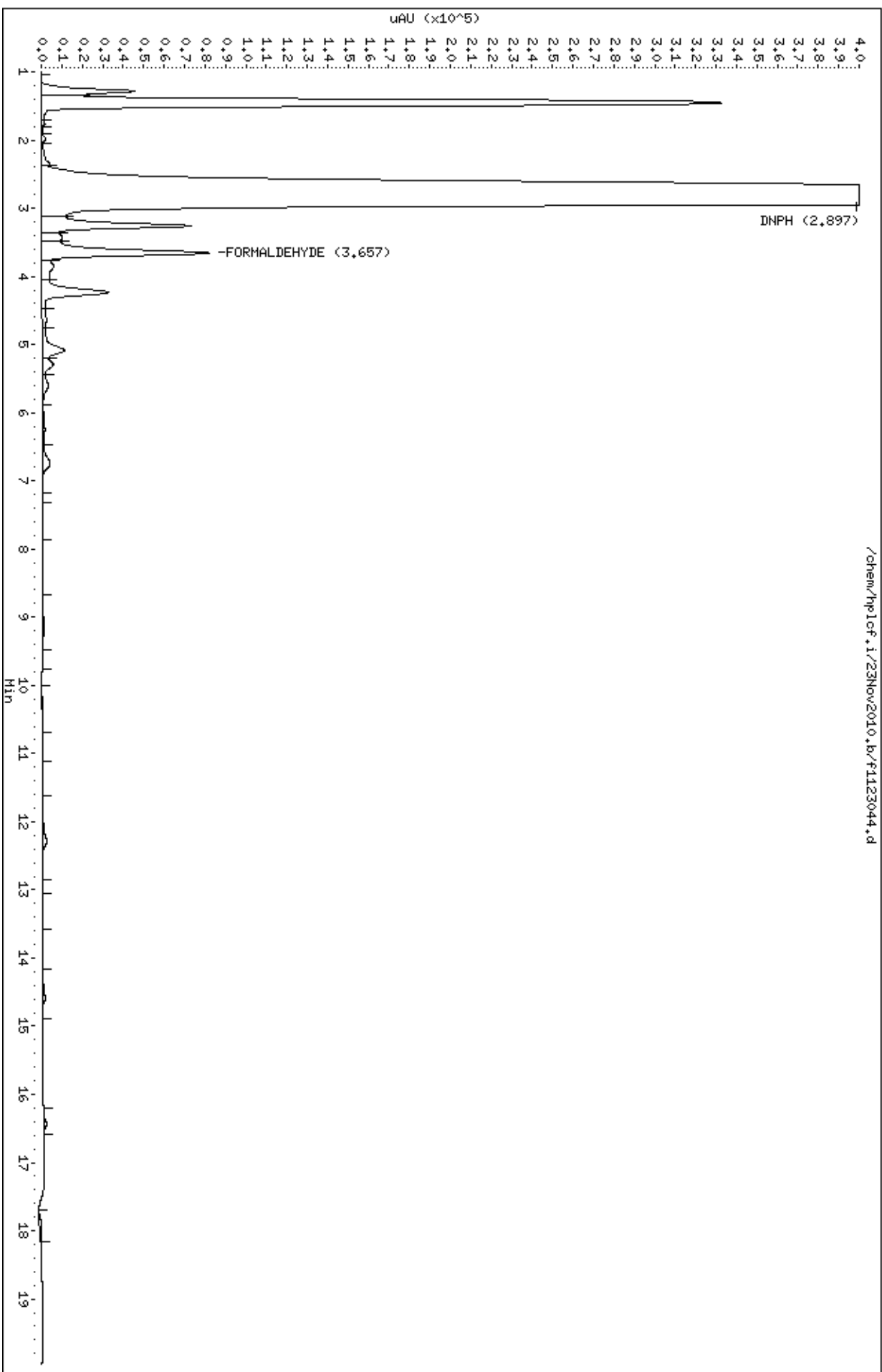
Sample Info: j1011418a-09a;

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



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

Client Sample ID: 118703

Lab ID#: 1011418A-10A

No Detections Were Found.

Client Sample ID: 118703

Lab ID#: 1011418A-10A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123036
Dil. Factor: 1.00

Date of Collection: NA
Date of Analysis: 11/23/10 07:51 PM
Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	1.0	0.52	Not Detected	Not Detected

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123036.d

Lab Smp Id: 1011418A-10A

Inj Date : 23-NOV-2010 19:51

Operator : FM

Smp Info : ;1011418A-10A;

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 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	RT	RESPONSE	ON-COLUMN	FINAL
	==	=====	=====		=====	(ng)	(ug)
1 DNPH	2.868	2.885	-0.017		33487410		
2 FORMALDEHYDE					Compound Not Detected.		

Data File: /chem/hp1cf.i/23Nov2010.b/f1123036.d

Page 1

Date : 23-NOV-2010 19:51

Client ID:

Instrument: hp1cf.i

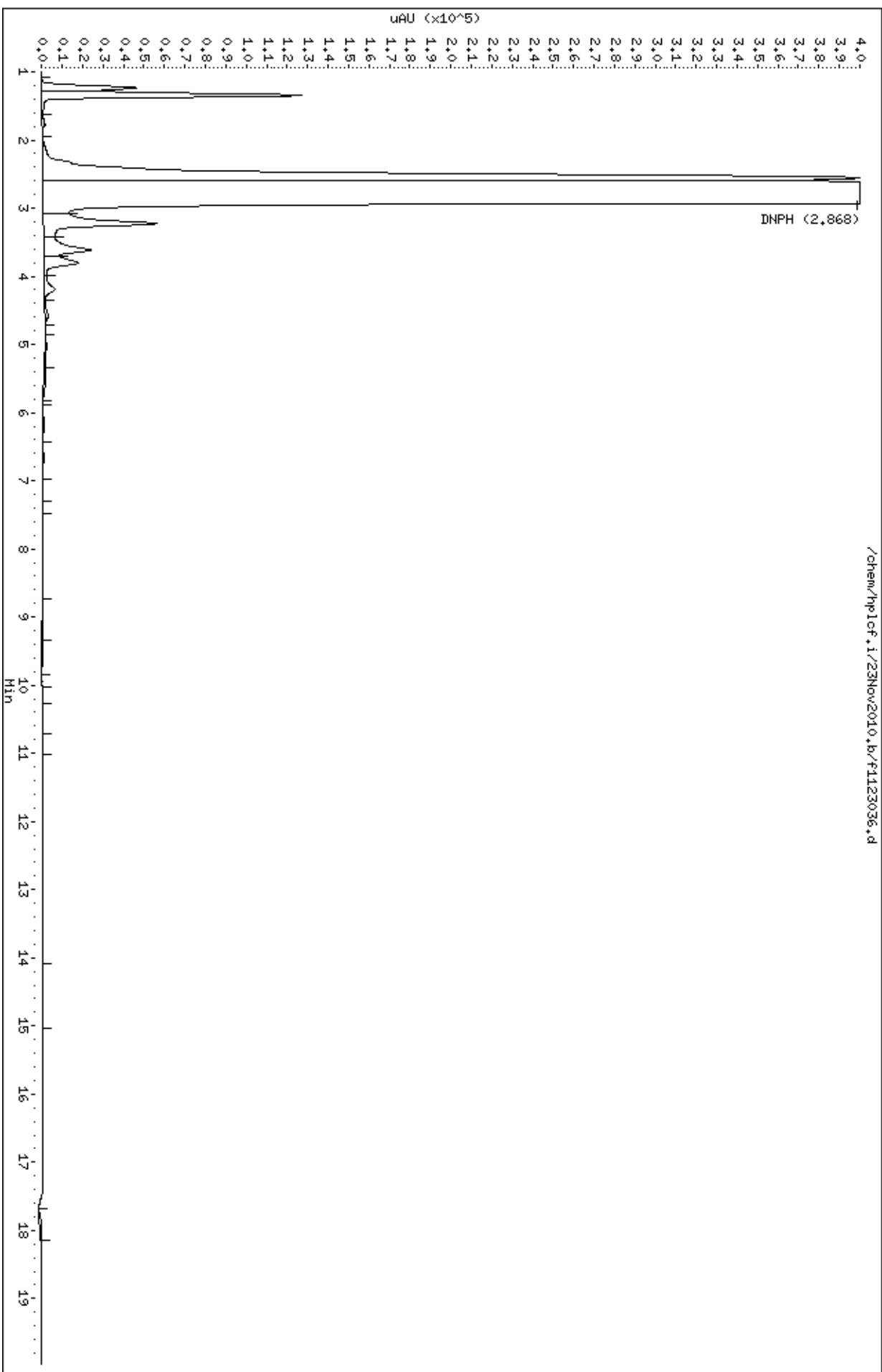
Sample Info: j1011418a-10a;

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118715

Lab ID#: 1011418A-11A

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	88	46

Client Sample ID: 118715

Lab ID#: 1011418A-11A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123057
Dil. Factor: 10.0

Date of Collection: 11/15/10
Date of Analysis: 11/24/10 03:10 AM
Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	88	46

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123057.d

Lab Smp Id: 1011418A-11AClient Smp ID: 10

Inj Date : 24-NOV-2010 03:10

Operator : FMInst ID: hplcf.i

Smp Info : ;1011418A-11A;10

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 10.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS					
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.868	2.885	-0.017	1584411		
2 FORMALDEHYDE	3.615	3.598	0.017	2407302	21.9897	88.0

Data File: /chem/hp1cf.i/23Nov2010.b/f1123057.d

Page 1

Date : 24-NOV-2010 03:10

Client ID: 10

Instrument: hp1cf.i

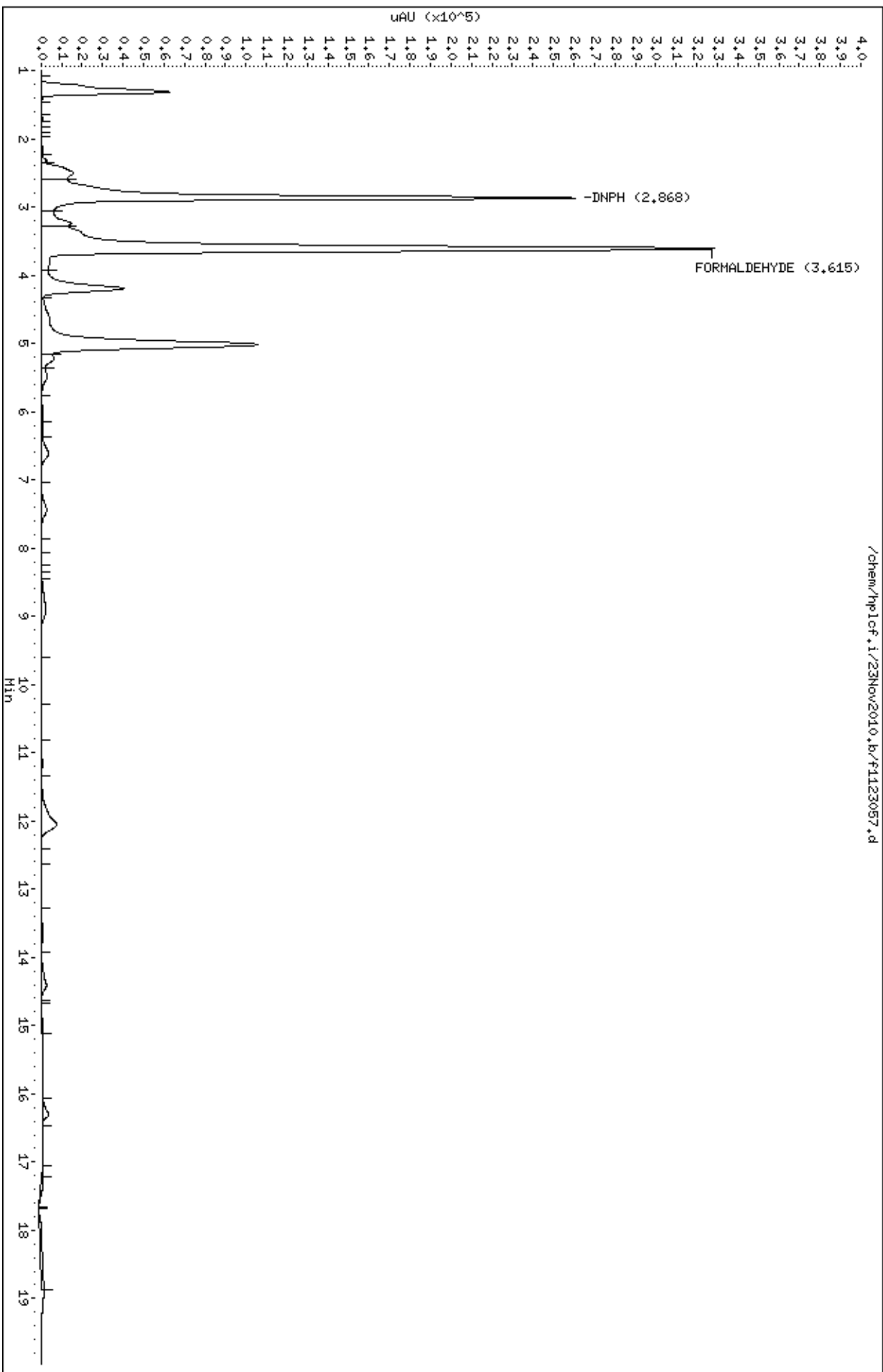
Sample Info: j10114180-11a;10

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118716

Lab ID#: 1011418A-12A

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	50	26

Client Sample ID: 118716

Lab ID#: 1011418A-12A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123058
Dil. Factor: 10.0

Date of Collection: 11/15/10
Date of Analysis: 11/24/10 03:30 AM
Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	50	26

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123058.d

Lab Smp Id: 1011418A-12AClient Smp ID: 10

Inj Date : 24-NOV-2010 03:30

Operator : FMInst ID: hplcf.i

Smp Info : ;1011418A-12A;10

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 10.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS					
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.788	2.885	-0.097	3227286		
2 FORMALDEHYDE	3.535	3.598	-0.063	1369798	12.5125	50.0

Data File: /chem/hp1cf.i/23Nov2010.b/f1123058.d

Page 1

Date : 24-NOV-2010 03:30

Client ID: 10

Instrument: hp1cf.i

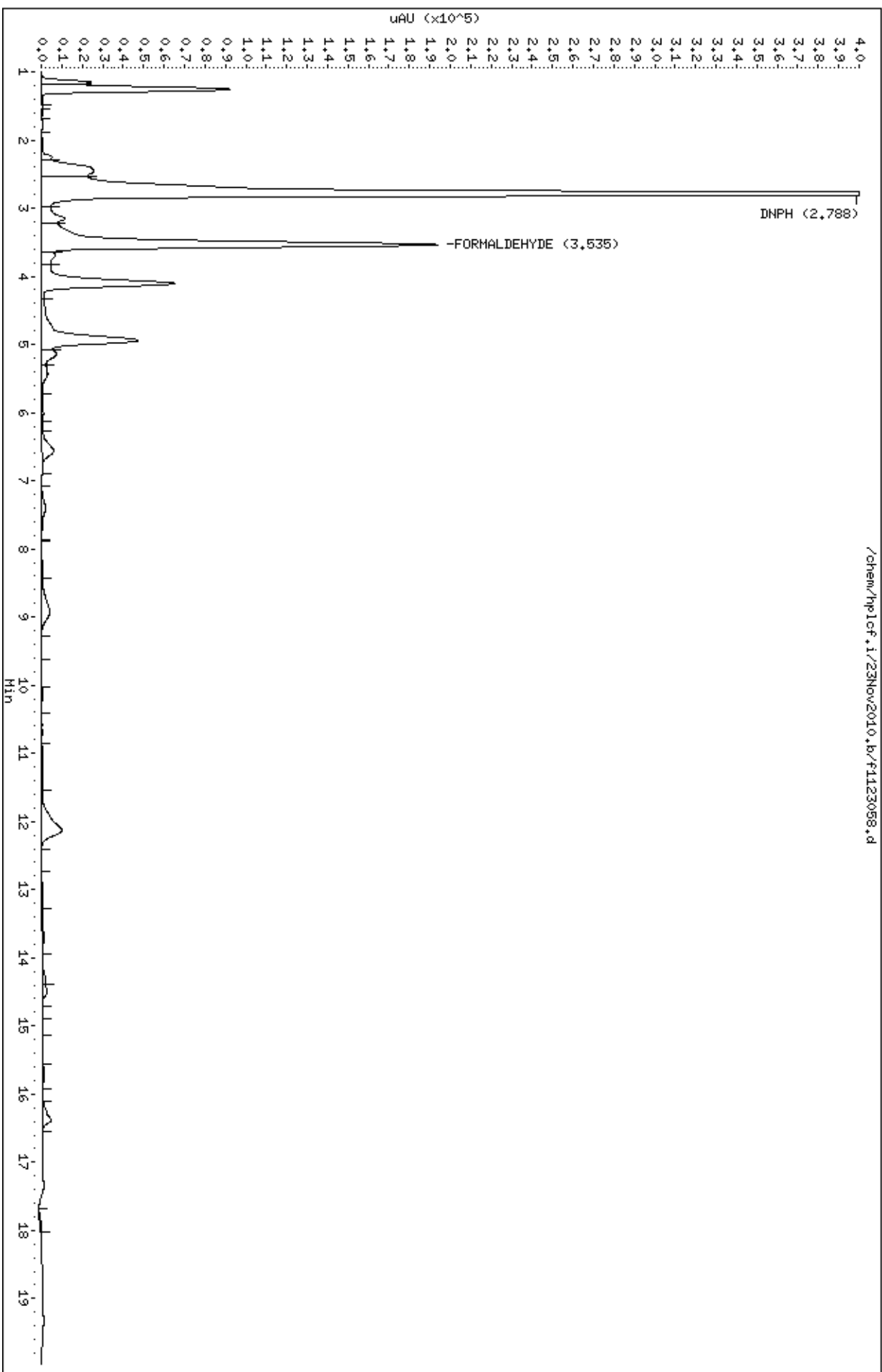
Sample Info: j10114180-120:10

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118717

Lab ID#: 1011418A-13A

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	71	36

Client Sample ID: 118717

Lab ID#: 1011418A-13A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123059
Dil. Factor: 10.0

Date of Collection: 11/15/10
Date of Analysis: 11/24/10 03:51 AM
Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	10	5.2	71	36

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,T0-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123059.d

Lab Smp Id: 1011418A-13AClient Smp ID: 10

Inj Date : 24-NOV-2010 03:51

Operator : FMInst ID: hplcf.i

Smp Info : ;1011418A-13A;10

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 10.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS					
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====	=====	=====	=====
1 DNPH	2.851	2.885	-0.034	2382924		
2 FORMALDEHYDE	3.604	3.598	0.006	1934594	17.6717	70.7

Data File: /chem/hp1cf.i/23Nov2010.b/f1123059.d

Page 1

Date : 24-NOV-2010 03:51

Client ID: 10

Instrument: hp1cf.i

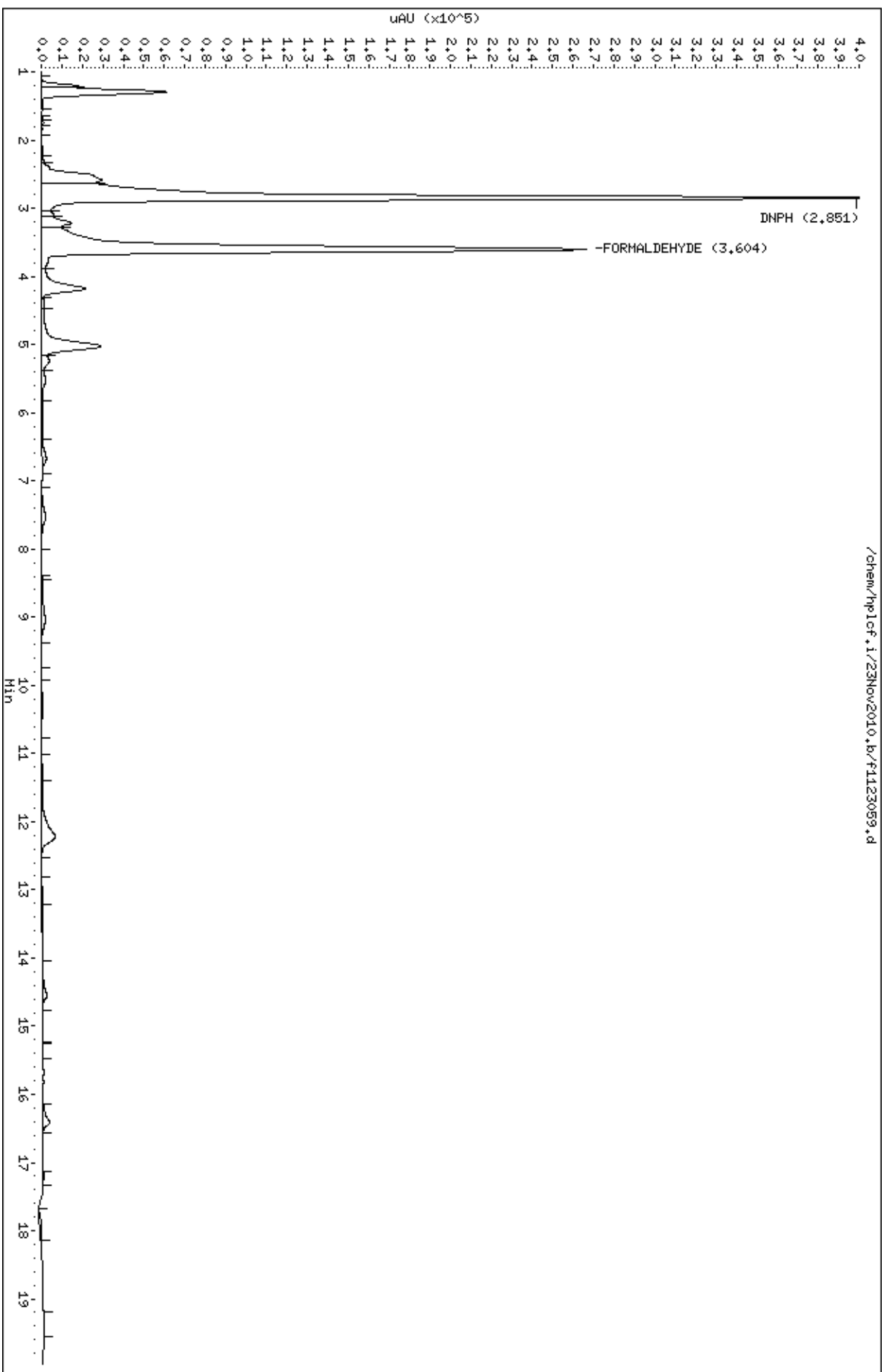
Sample Info: #10114180-13A:10

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Summary of Detected Compounds
RADIELLO ANALYSIS BY HPLC/UV

Client Sample ID: 118718

Lab ID#: 1011418A-14A

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	1.0	0.52	2.5	1.3

Client Sample ID: 118718

Lab ID#: 1011418A-14A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123045

Date of Collection: 11/15/10

Dil. Factor: 1.00

Date of Analysis: 11/23/10 10:59 PM

Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	1.0	0.52	2.5	1.3

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123045.d

Lab Smp Id: 1011418A-14A

Inj Date : 23-NOV-2010 22:59

Operator : FM

Smp Info : ;1011418A-14A;

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 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	2.869	2.885	-0.016	18999471		
2 FORMALDEHYDE	3.629	3.598	0.031	689184	6.29540	2.52

Data File: /chem/hp1cf.i/23Nov2010.b/f1123045.d

Page 1

Date : 23-NOV-2010 22:59

Client ID:

Instrument: hp1cf.i

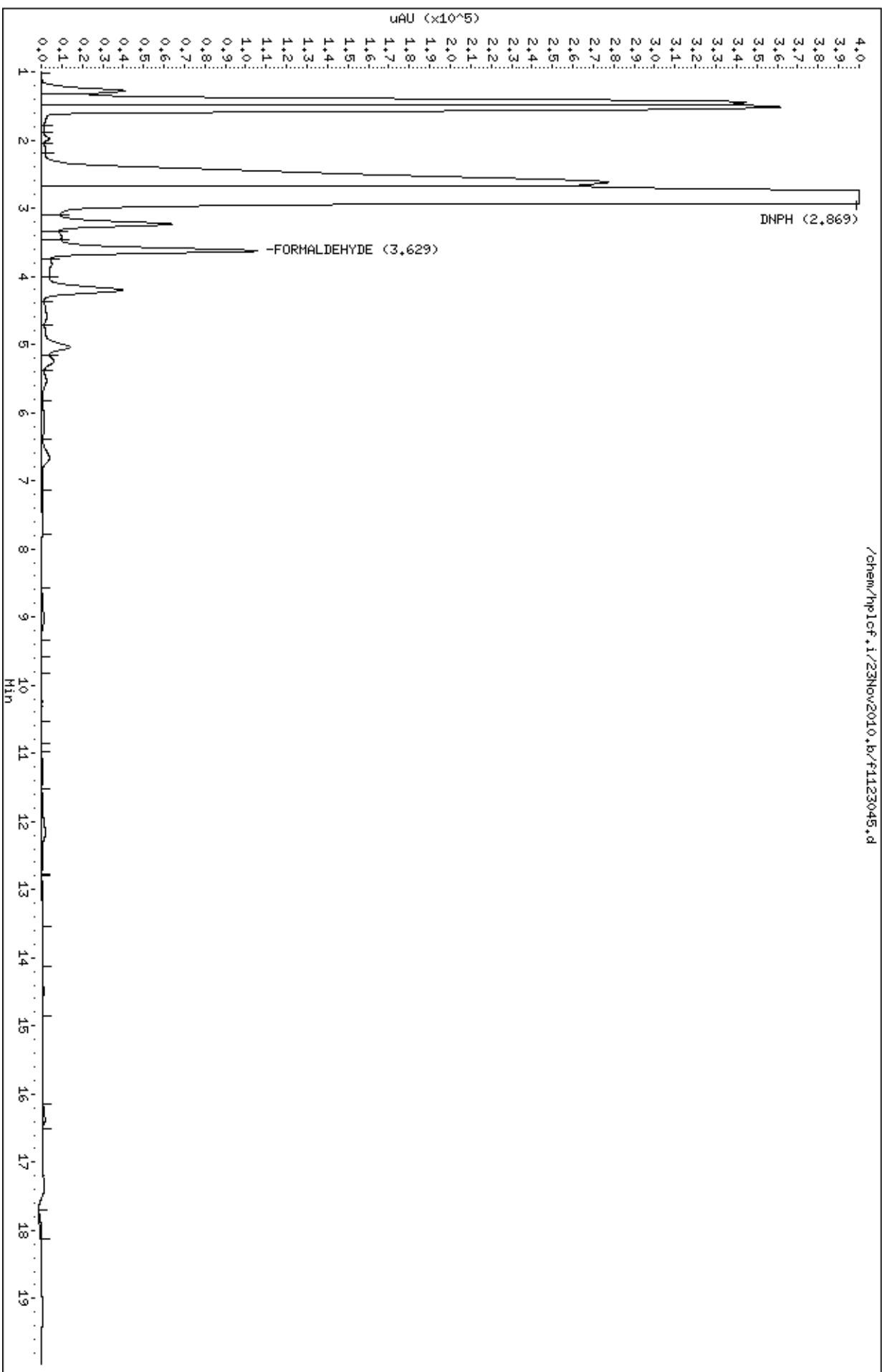
Sample Info: j10114180-14a;

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



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

Client Sample ID: 118719

Lab ID#: 1011418A-15A

No Detections Were Found.

Client Sample ID: 118719

Lab ID#: 1011418A-15A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123037
Dil. Factor: 1.00

Date of Collection: 11/15/10
Date of Analysis: 11/23/10 08:12 PM
Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	1.0	0.52	Not Detected	Not Detected

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123037.d

Lab Smp Id: 1011418A-15A

Inj Date : 23-NOV-2010 20:12

Operator : FM

Smp Info : ;1011418A-15A;

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 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	2.882	2.885	-0.003	25837014		
2 FORMALDEHYDE				Compound Not Detected.		

Data File: /chem/hplcf.i/23Nov2010.b/f1123037.d

Page 1

Date : 23-NOV-2010 20:12

Client ID:

Instrument: hplcf.i

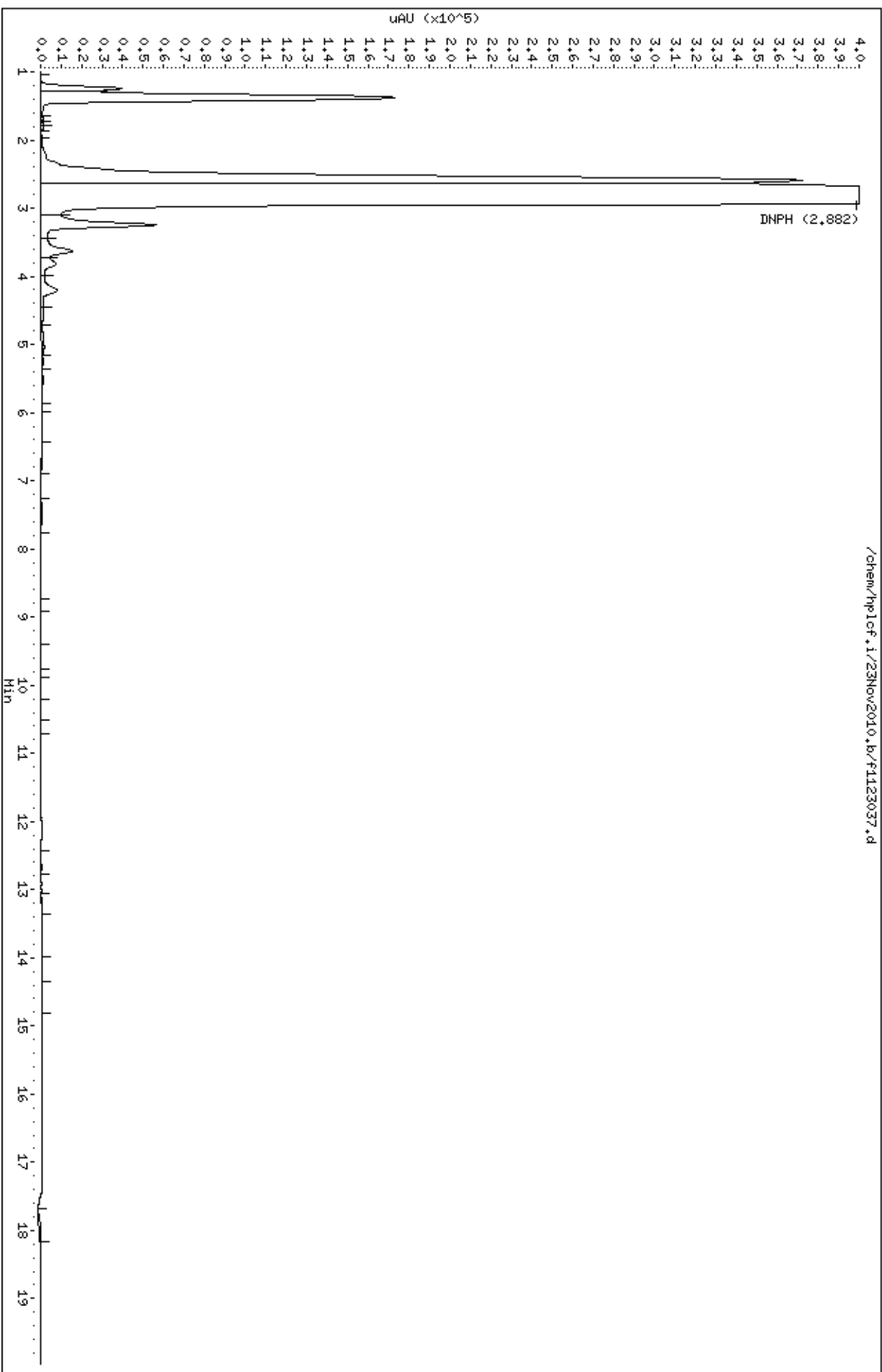
Sample Info: j10114180-15A;

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



QC Results and Raw Data

Client Sample ID: Lab Blank

Lab ID#: 1011418A-16A

RADIELLO ANALYSIS BY HPLC/UV

File Name:	f1123010	Date of Collection: NA
Dil. Factor:	1.00	Date of Analysis: 11/23/10 10:48 AM
		Date of Extraction: 11/23/10

Compound	Rpt. Limit (ug)	Rpt. Limit (ug/m3)	Amount (ug)	Amount (ug/m3)
Formaldehyde	1.0	0.52	Not Detected	Not Detected

Container Type: Radiello 165 (Aldehydes)

Data File: /chem/hplcf.i/23Nov2010.b/f1123010.d

Page 1

Date : 23-NOV-2010 10:48

Client ID: Lab Blank

Instrument: hplcf.i

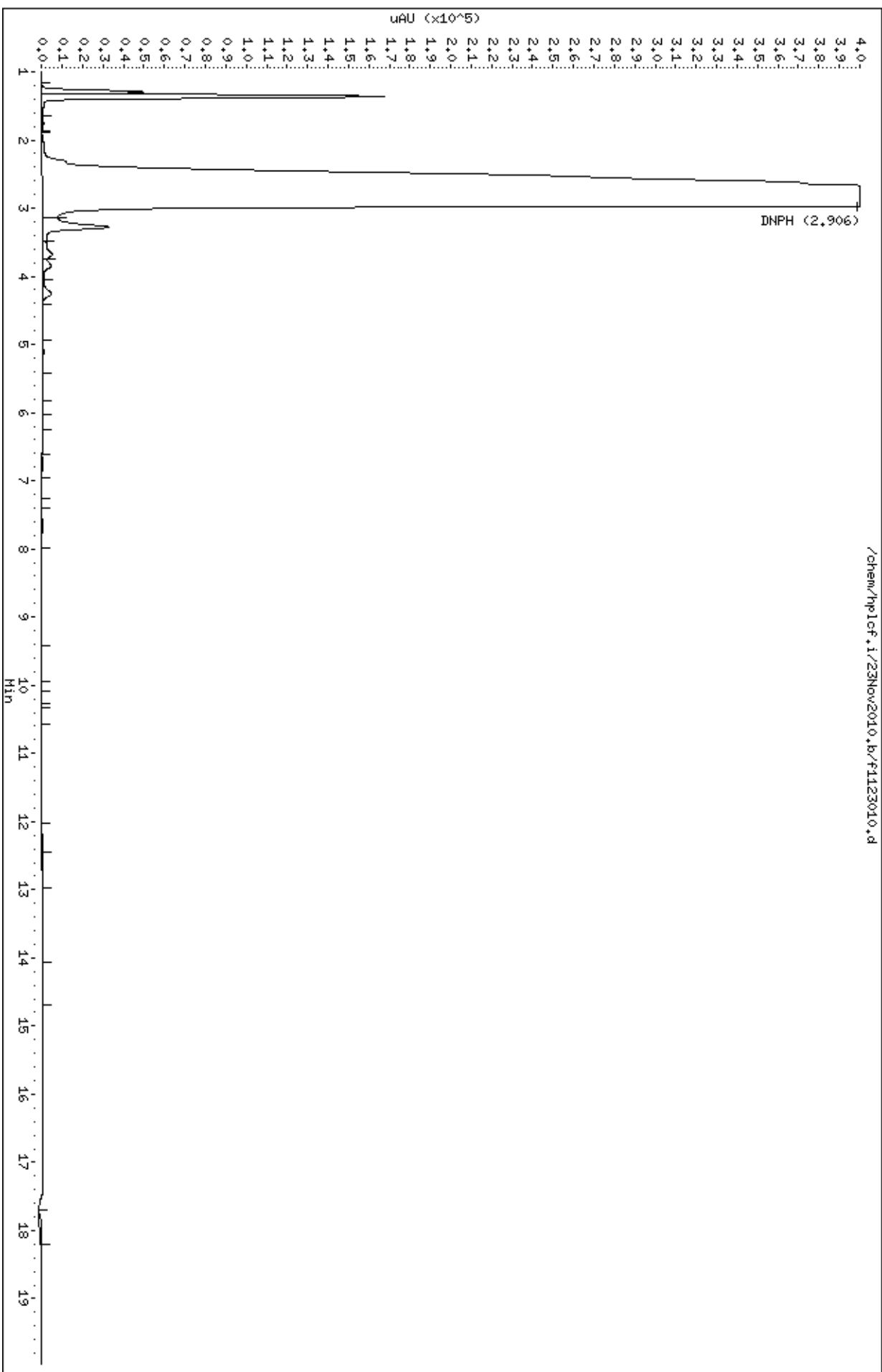
Sample Info: f10114180;Lab Blank

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



SAMPLE RESULTS/SAMPLE RESULTS DUPLICATE

Lab Name: Air Toxics Ltd.

Lab File ID: f1123020.d & f1123019.d

Lab Sample ID: &

Dilution: 1.00 & 1.00

Client Sample ID: LCS & LCSD

Date Analyzed: 11/23/10 & 11/23/10

CAS Number	Compound	Original		Duplicate		RPD	Result Less Than 5X RL
		Amount	Flags	Amount	Flags		
50-00-0	Formaldehyde	99		99		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

	0.05000	0.10000	0.25000	0.50000	1.000	4.000	—	
Compound	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6	RRF	% RSD
	-----	-----	-----	-----	-----	-----		
	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
-----	-----	-----	-----	-----	-----	-----	-----	-----

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

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 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	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.dCompound: CotanolInitials: CRL

Sample 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

Cary Lee
Signed

9/15/10
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

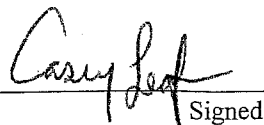
Calculation Check:

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: _____


Signed9/15/10
Date

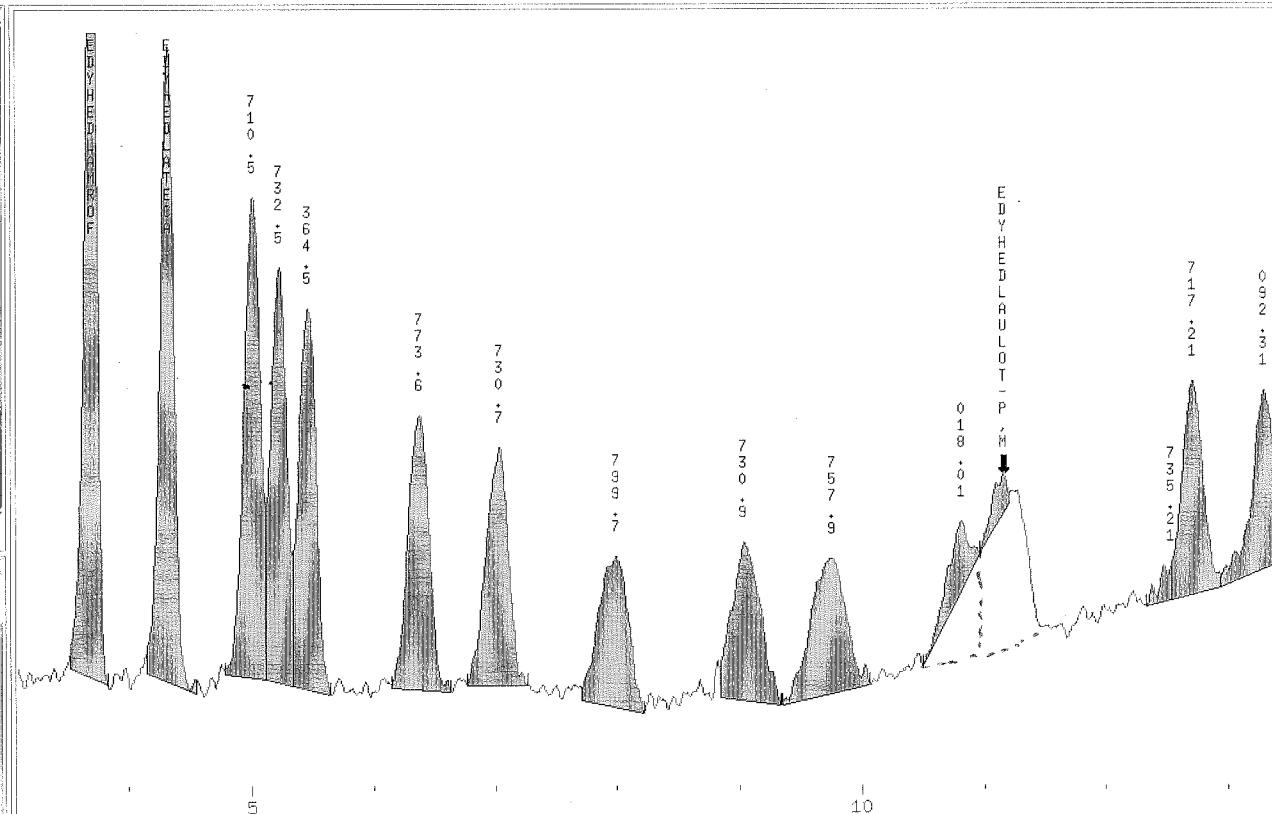
Revised 07/08

File Record Edit 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

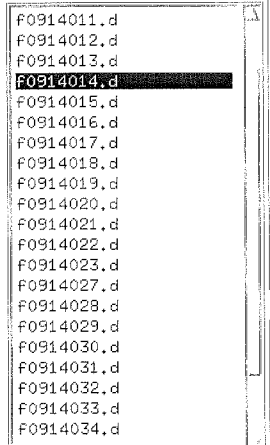
1 DNPH
+ 2 FORMALDEHYDE
+ 3 ACETALDEHYDE
* 15 M,P-TOLUALDEHY

Compound: M,P-TOLUAL

Update Method Expected
Time New Time

11.157

Done Apply Unsa



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

- Mark M,P-TOLUALDEHYDE Undetected.

09/15/10

9/15/10

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/14Sep2010.b/f0914040.d

Lab Smp Id: 1838-36-4.0Client Smp ID: LCS

Inj Date : 15-SEP-2010 03:33

Operator : CRLInst 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 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1QC Sample: LCS

Dil Factor: 1.00000

Integrator: FalconCompound Sublist: all.sub

Target Version: 3.50

Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve 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

Page 1

Date : 15-SEP-2010 03:33

Client ID: LCS

Instrument: hp1cf.i

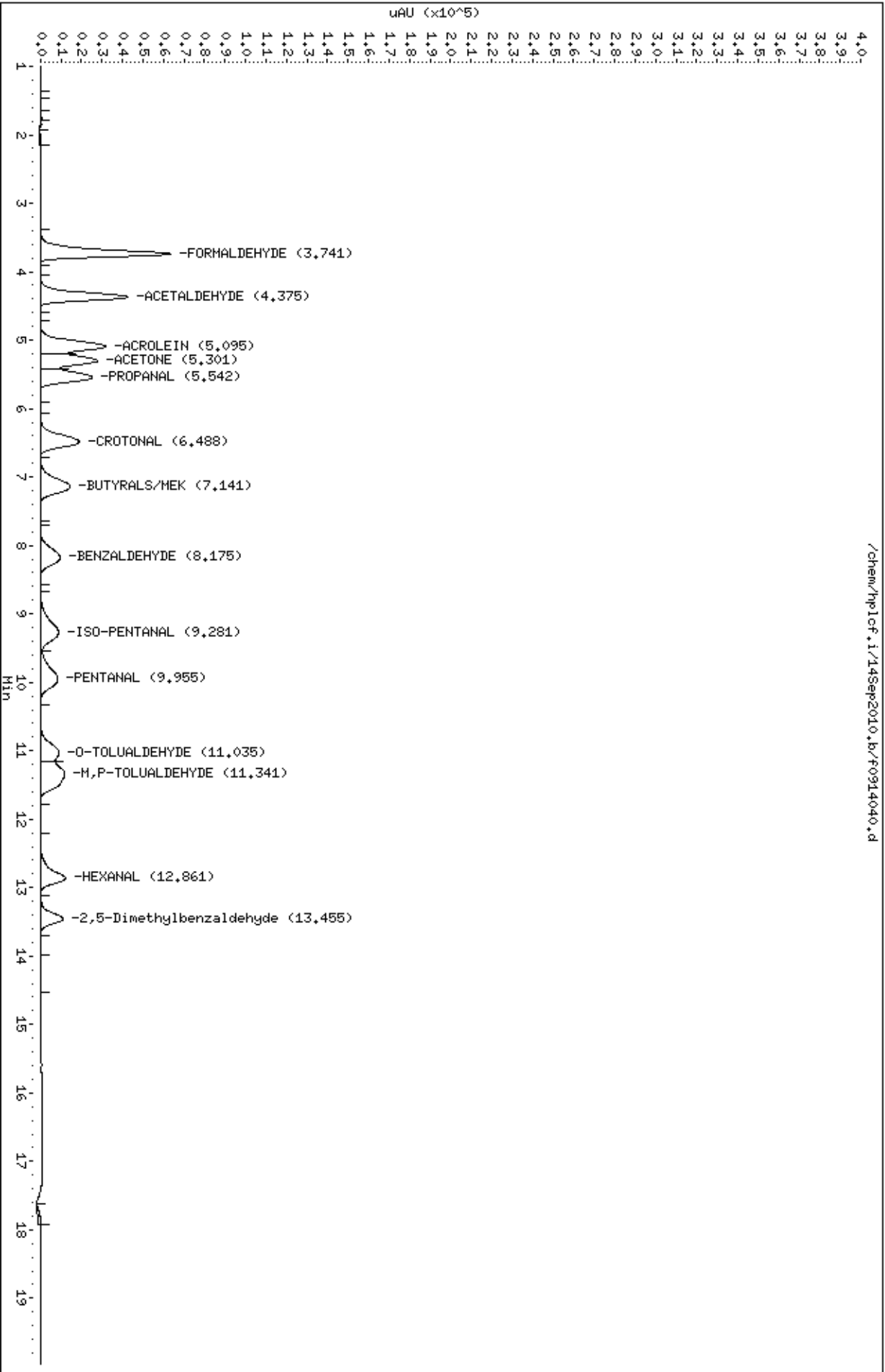
Sample Info: f1838-36-4.0;LCS

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

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

Lab Smp Id: 1838-34-0.05Client Smp ID: 1

Inj Date : 14-SEP-2010 17:27

Operator : CRLInst 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 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 17:27Cal File: f0914011.d

Als bottle: 1Calibration Sample, Level: 1

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

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

Data File: /chem/hp1cf.i/14Sep2010.b/f0914011.d

Page 1

Date : 14-SEP-2010 17:27

Client ID: 1

Instrument: hp1cf.i

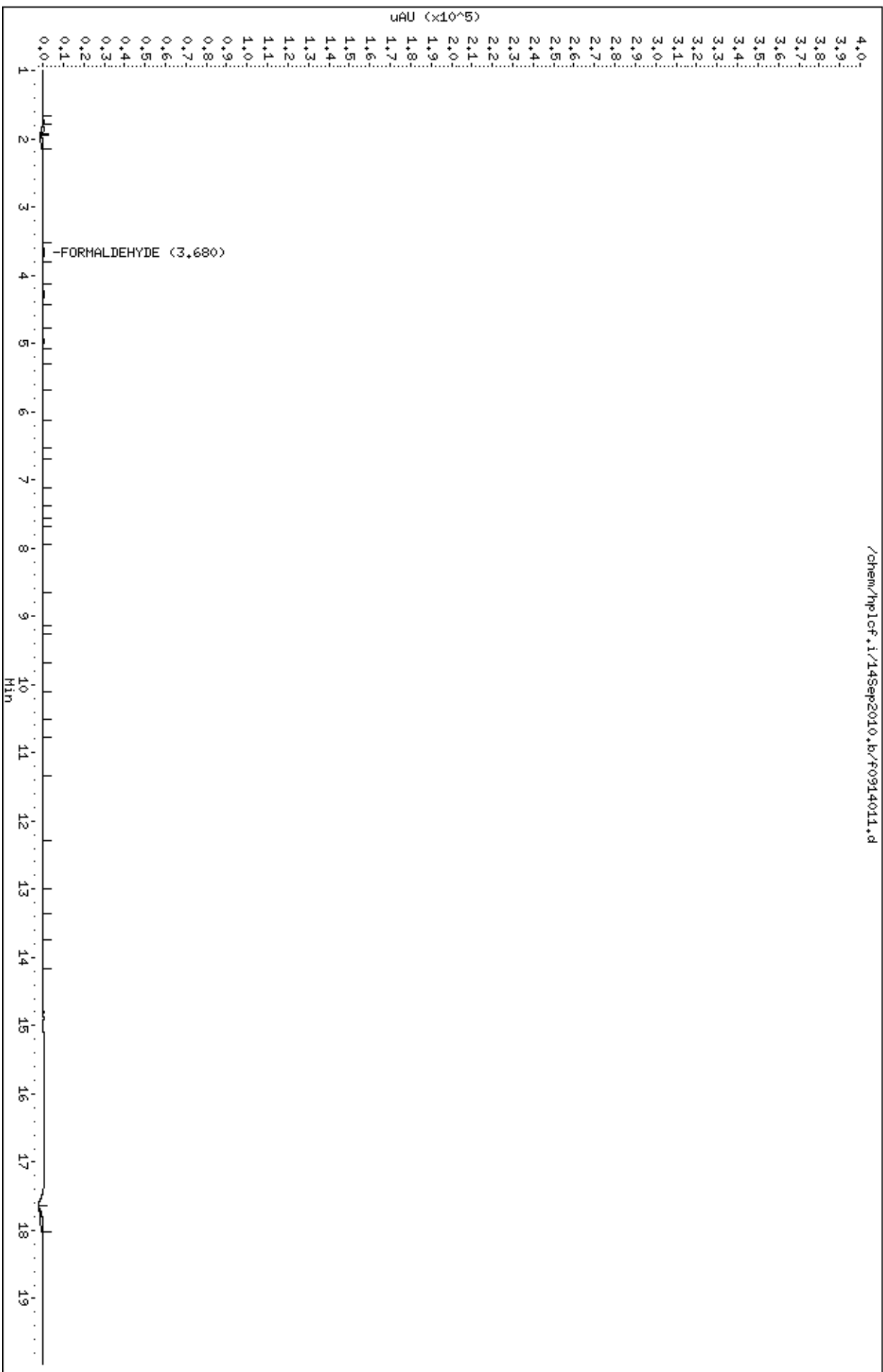
Sample Info: f1838-34-0.05;1

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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	AMOUNTS					
	CAL-AMT			ON-COL		
	RT	EXP RT	DLT RT	RESPONSE	(UG)	(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

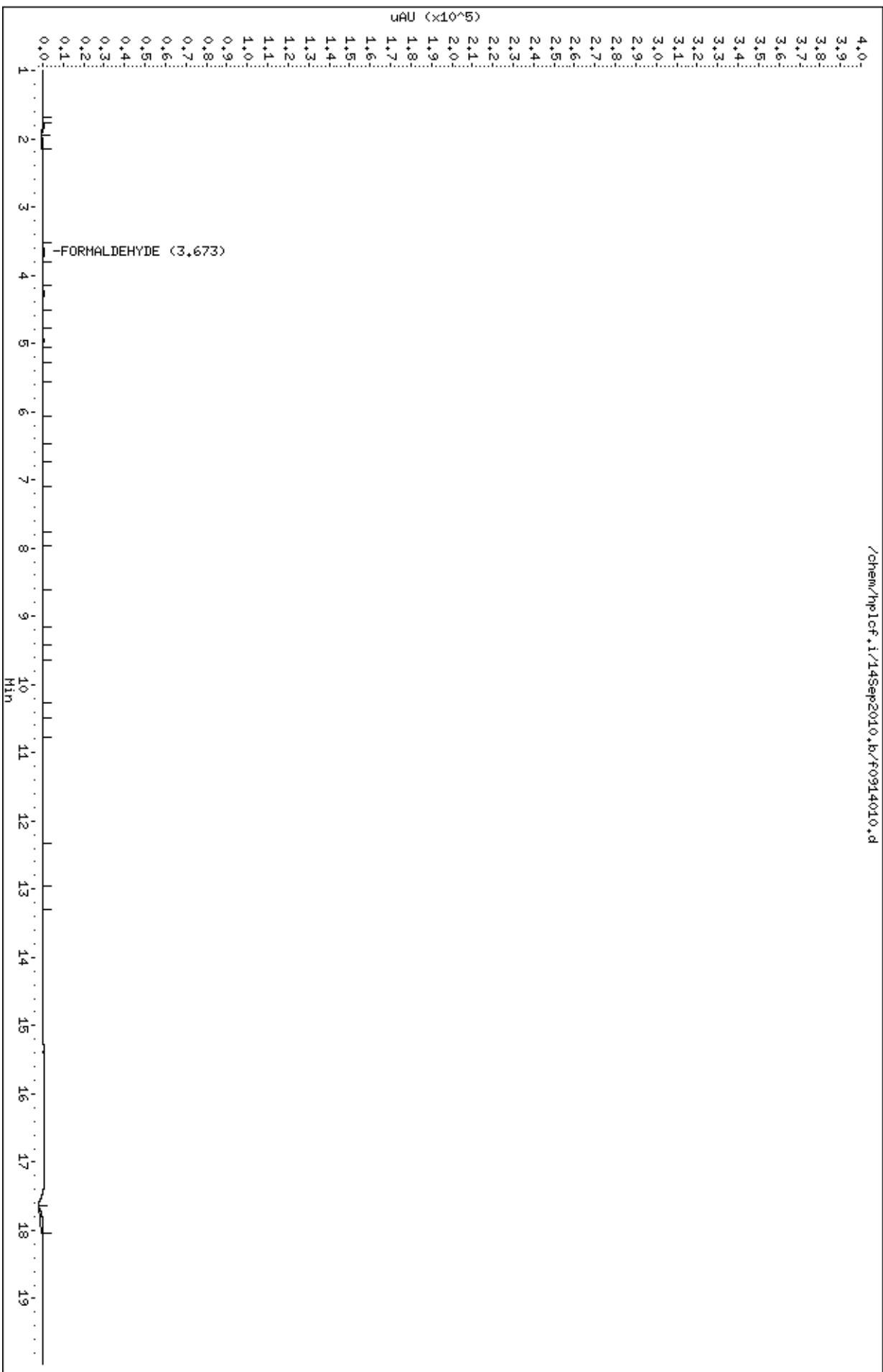
Sample Info: #1838-34-0.05;1

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/14Sep2010.b/f0914009.d

Lab Smp Id: 1838-34-0.05Client Smp ID: 1

Inj Date : 14-SEP-2010 16:46

Operator : CRLInst 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 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 16:46Cal File: f0914009.d

Als bottle: 1Calibration Sample, Level: 1

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	AMOUNTS					
	CAL-AMT			ON-COL		
	(UG)			(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

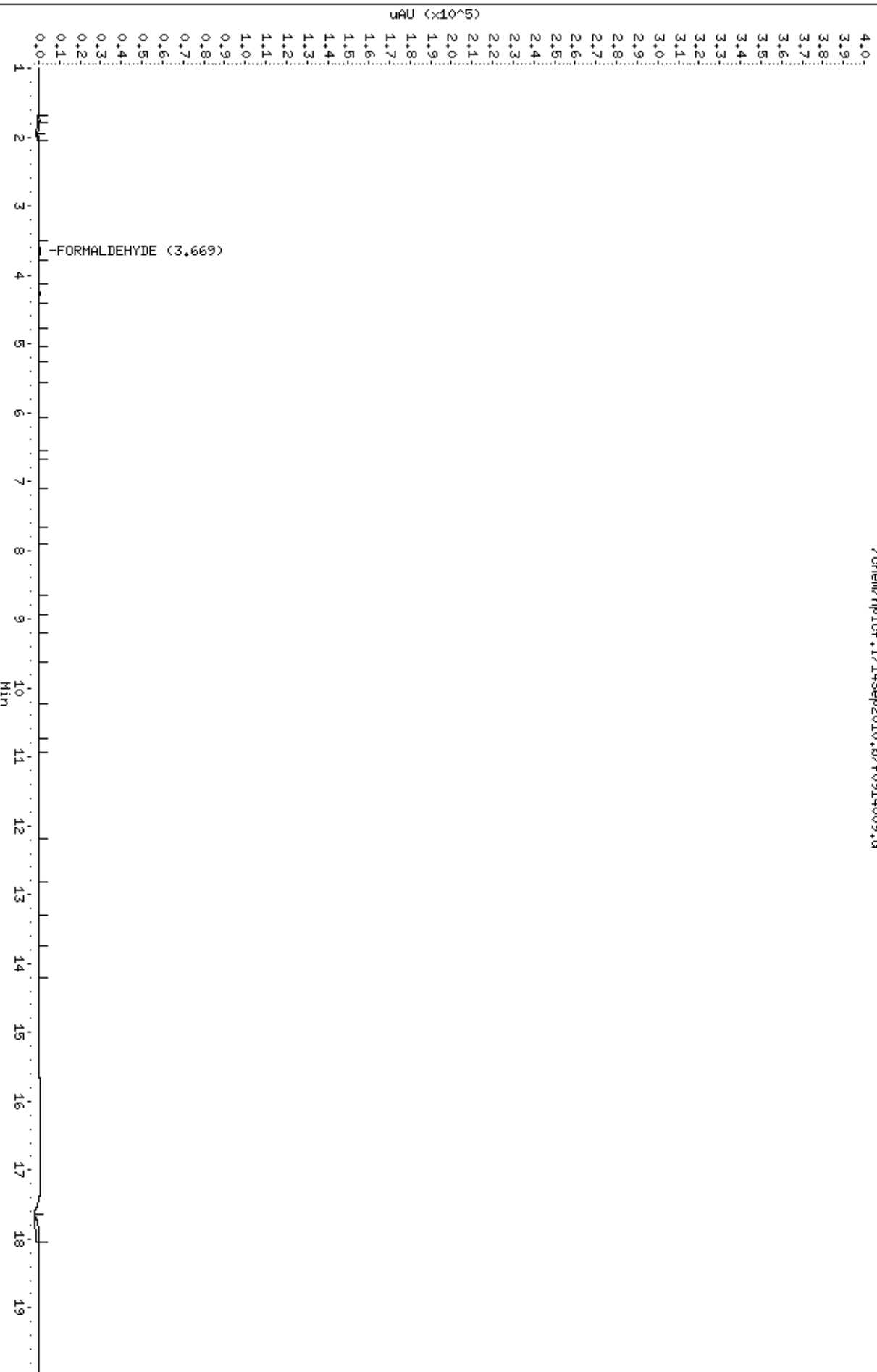
Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00

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



Aldehydes/Ketones To-11, TO-5, CARB430, 0011, Radiello 165

Cpnd Variable	Local Compound Variable
1	1
2	2
3	3
4	4
5	5
6	6
7	7
8	8
9	9
10	10
11	11
12	12
13	13
14	14
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69	69
70	70
71	71
72	72
73	73
74	74
75	75
76	76
77	77
78	78
79	79
80	80
81	81
82	82
83	83
84	84
85	85
86	86
87	87
88	88
89	89
90	90
91	91
92	92
93	93
94	94
95	95
96	96
97	97
98	98
99	99
100	100

M - Compound response manually integrated.

Data File: /chem/hp1cf.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

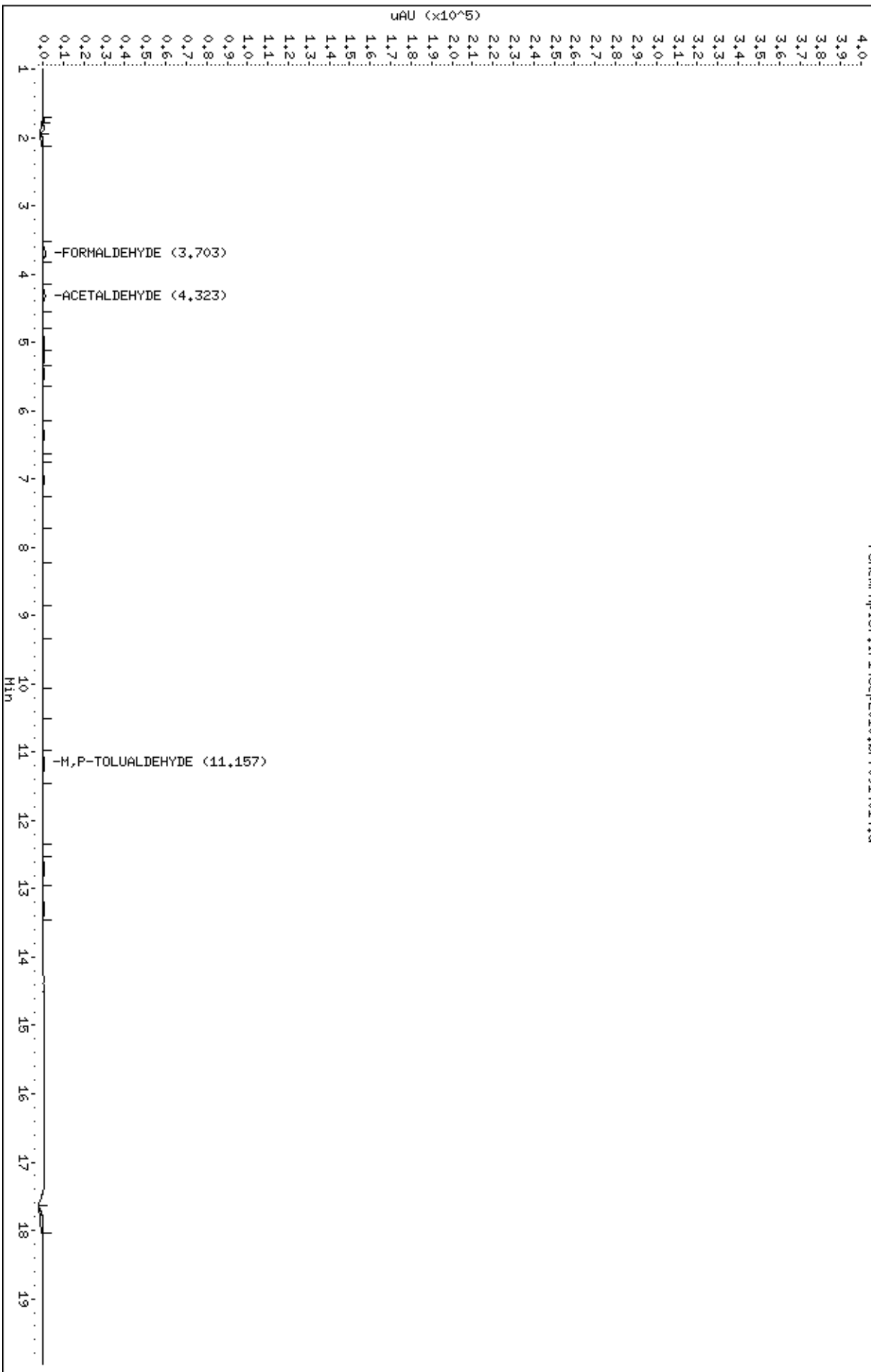
Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00

/chem/hp1cf.i/14Sep2010.b/f0914014.d



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds						AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE		CAL-AMT	ON-COL
	==	=====	=====	=====		(UG)	(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

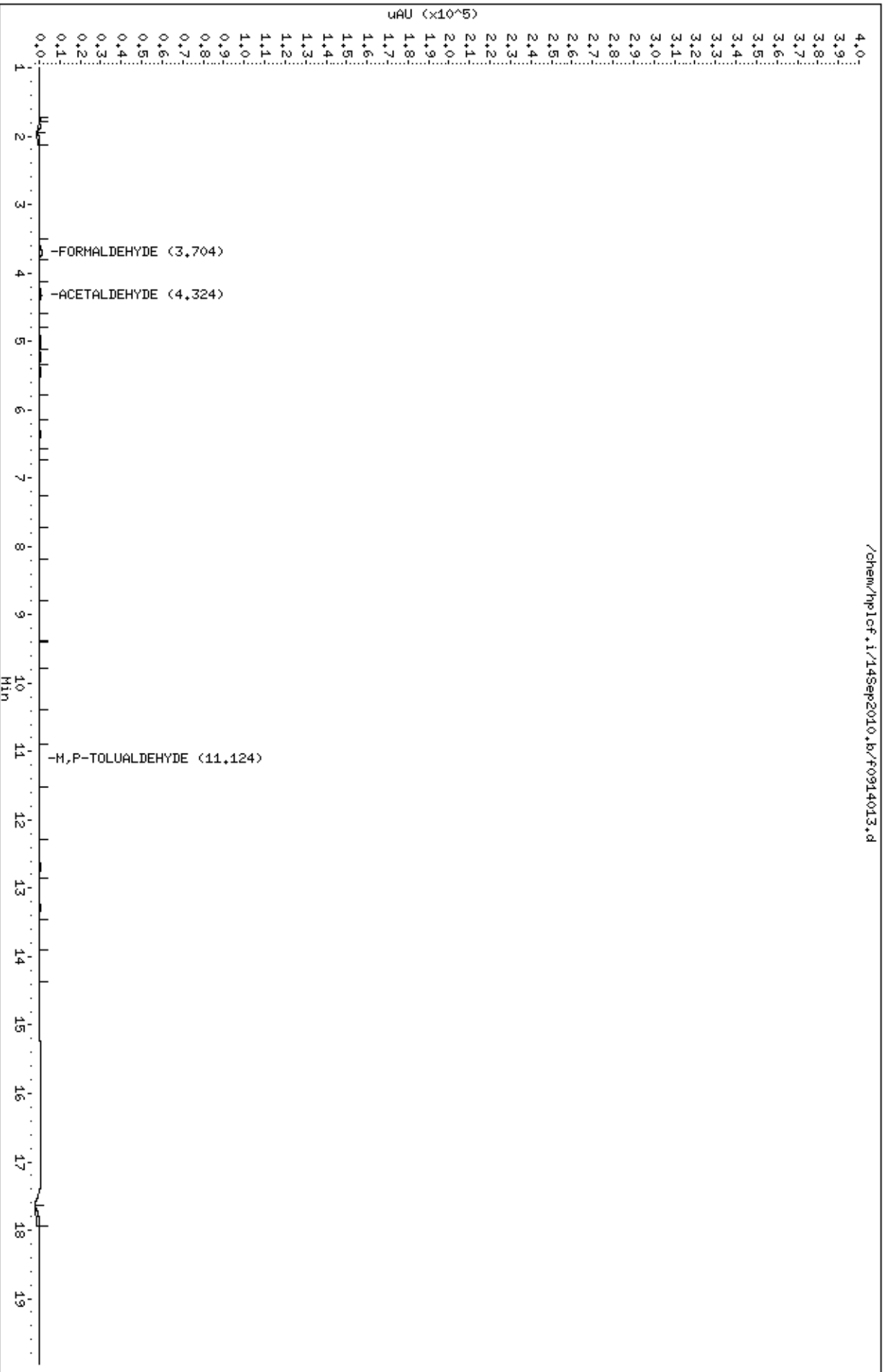
Sample Info: f1838-34-0.1;2

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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

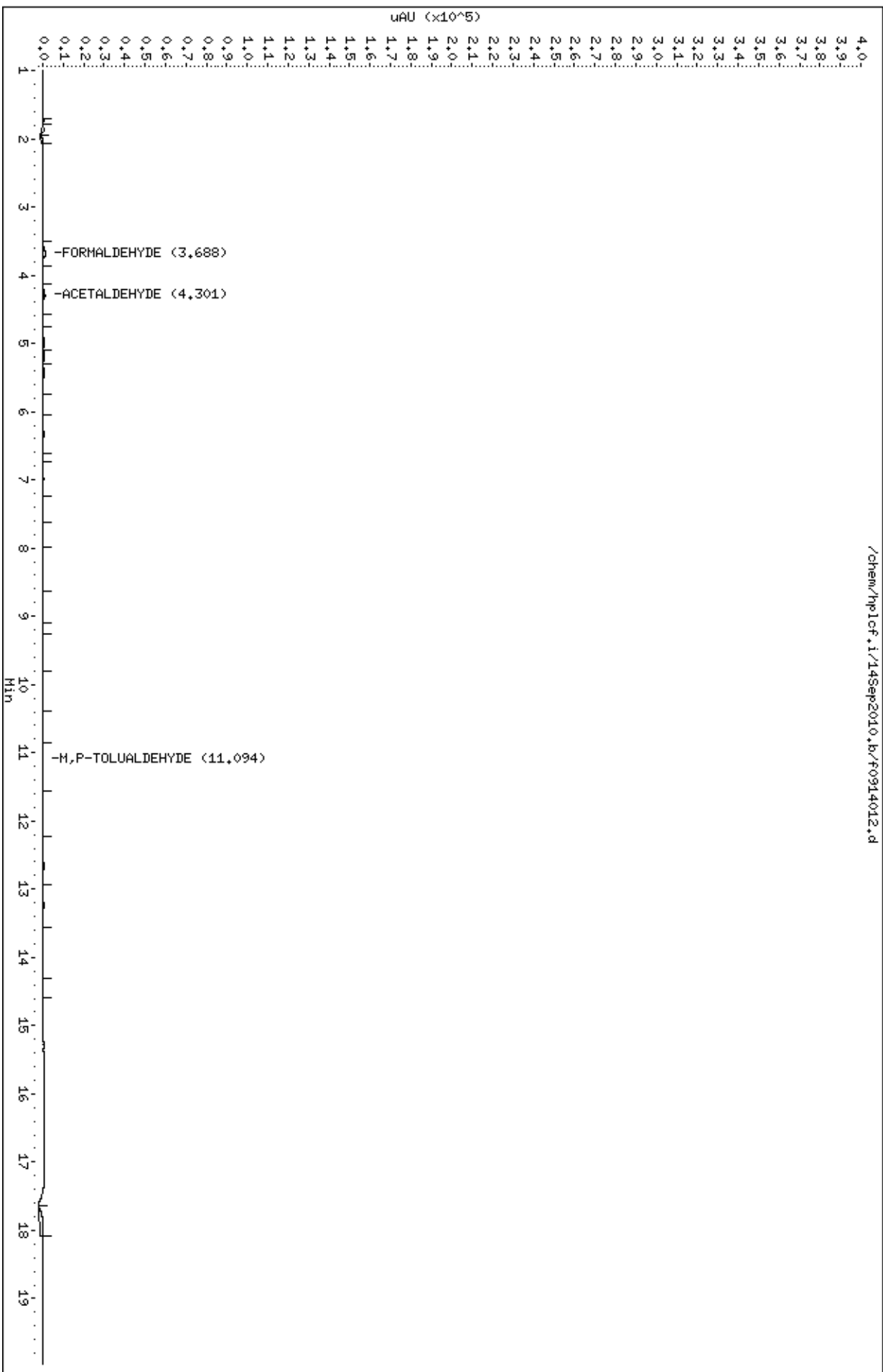
Sample Info: f1838-34-0.1;2

Purge Volume: 5.0

Operator: CRL

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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 ACROLEIN	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: #1838-34-0.25;3

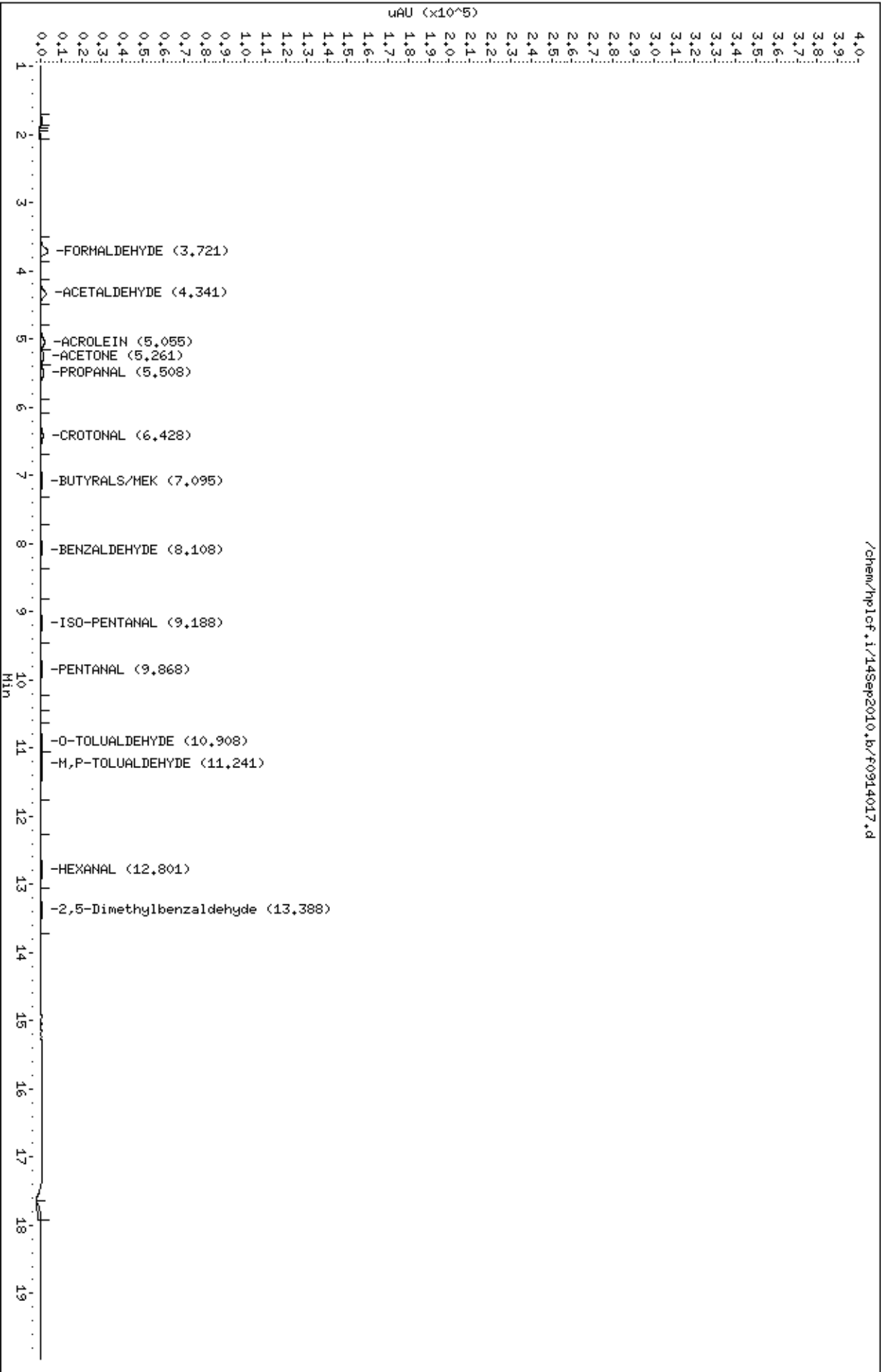
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve 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 ACROLEIN	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

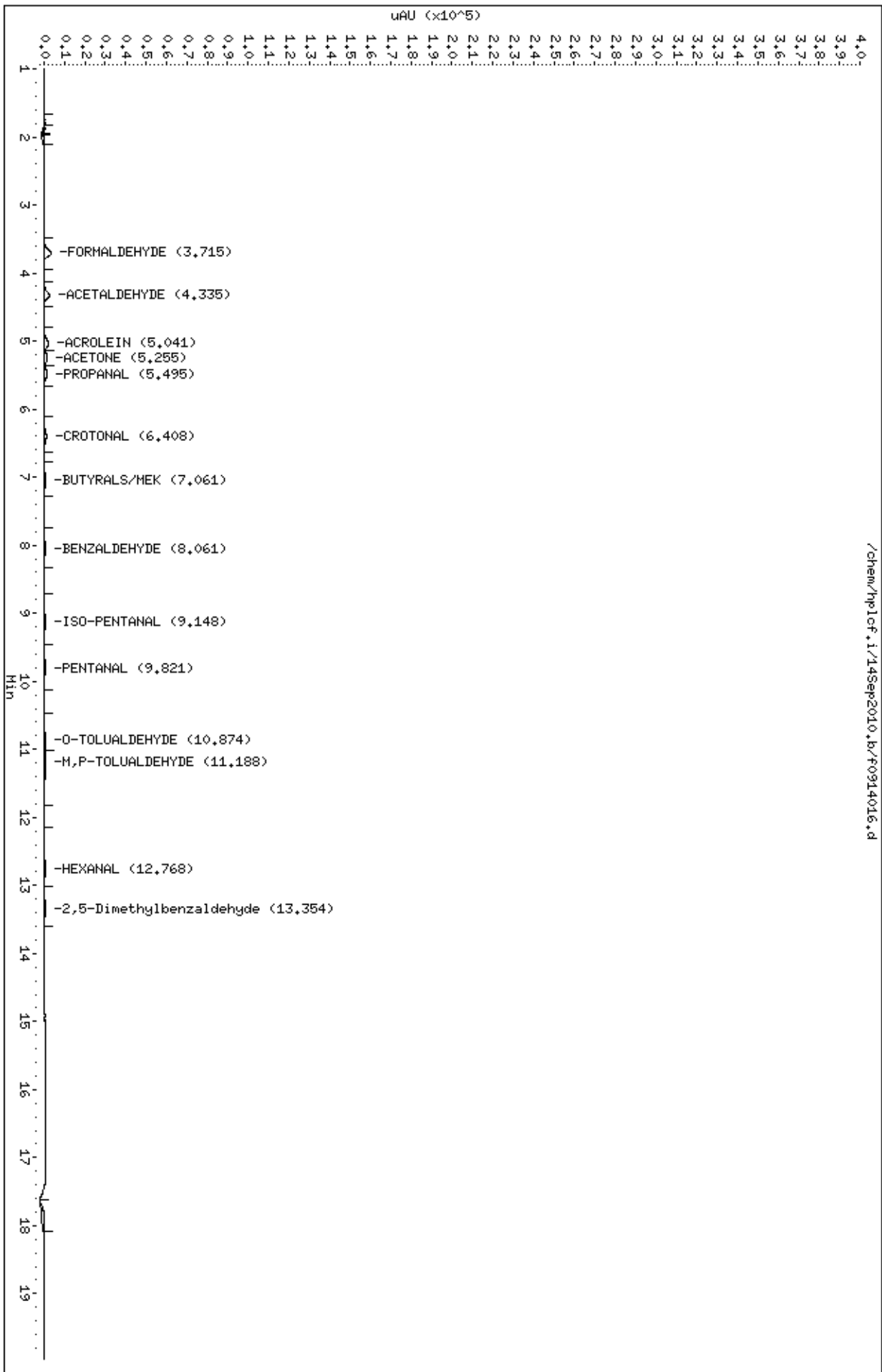
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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

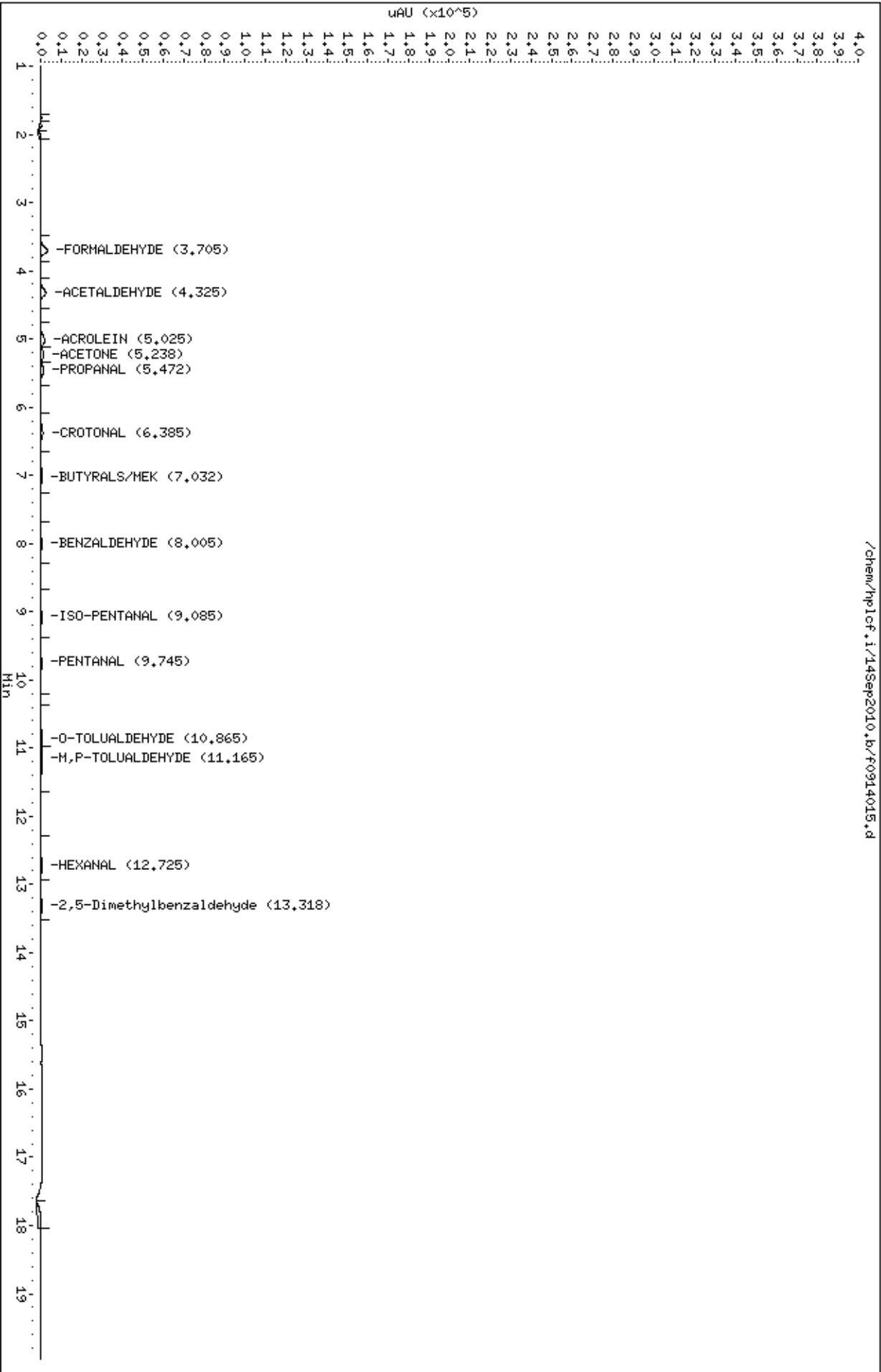
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



/chem/hp1cf.i/14Sep2010.b/f0914015.d

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve 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 ACROLEIN	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

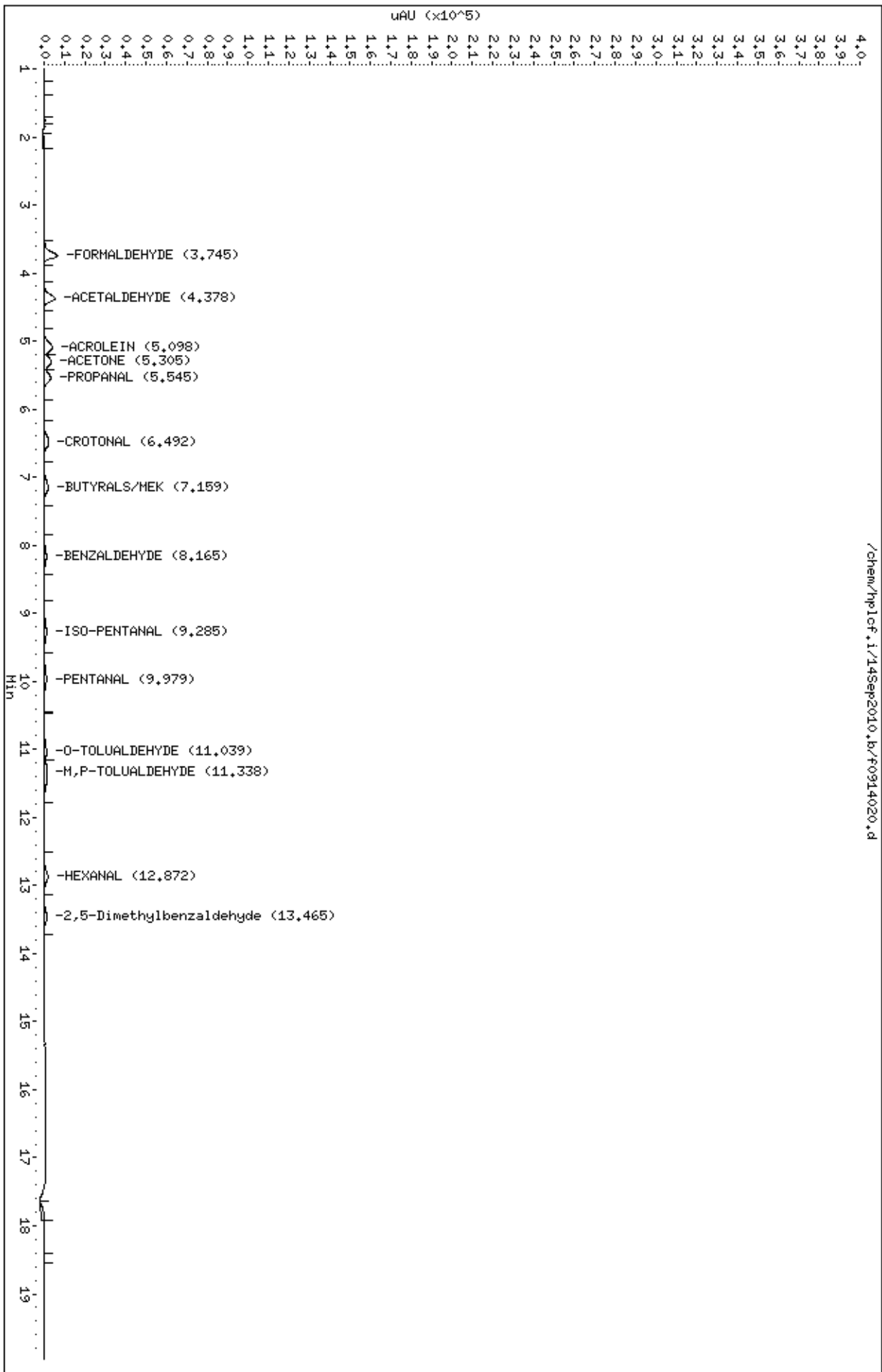
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/14Sep2010.b/f0914019.d

Lab Smp Id: 1838-34-0.5Client Smp ID: 4

Inj Date : 14-SEP-2010 20:15

Operator : CRLInst 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 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 20:15Cal File: f0914019.d

Als bottle: 1Calibration Sample, Level: 4

Dil Factor: 1.00000

Integrator: FalconCompound Sublist: all.sub

Target Version: 3.50

Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds						AMOUNTS	
	RT	EXP RT	DLT RT	RT	RESPONSE	CAL - AMT	ON - COL
						(UG)	(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: J1838-34-0.5;4

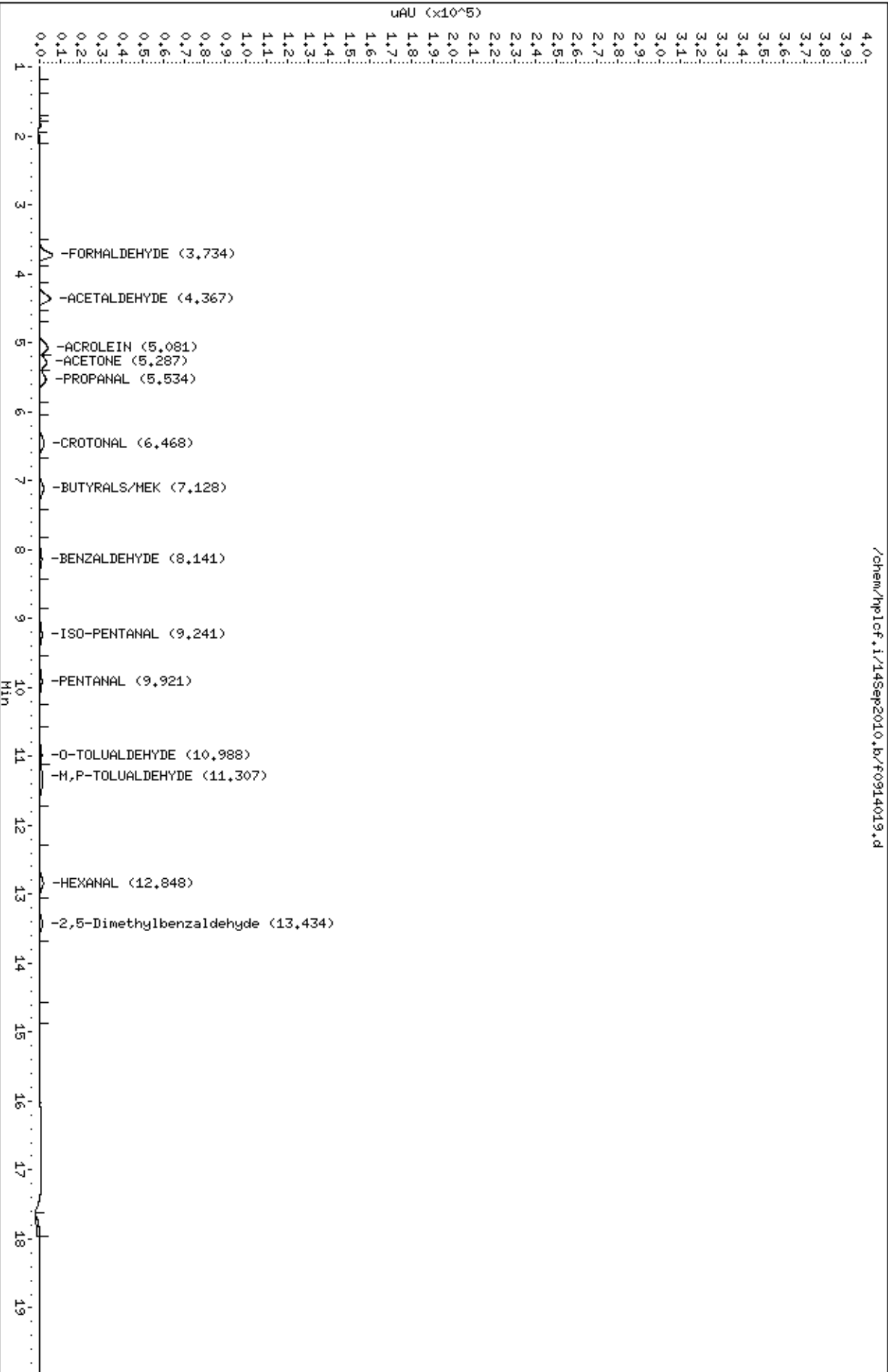
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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	RT	RESPONSE	CAL - AMT	ON - COL
						(UG)	(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 ACROLEIN	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

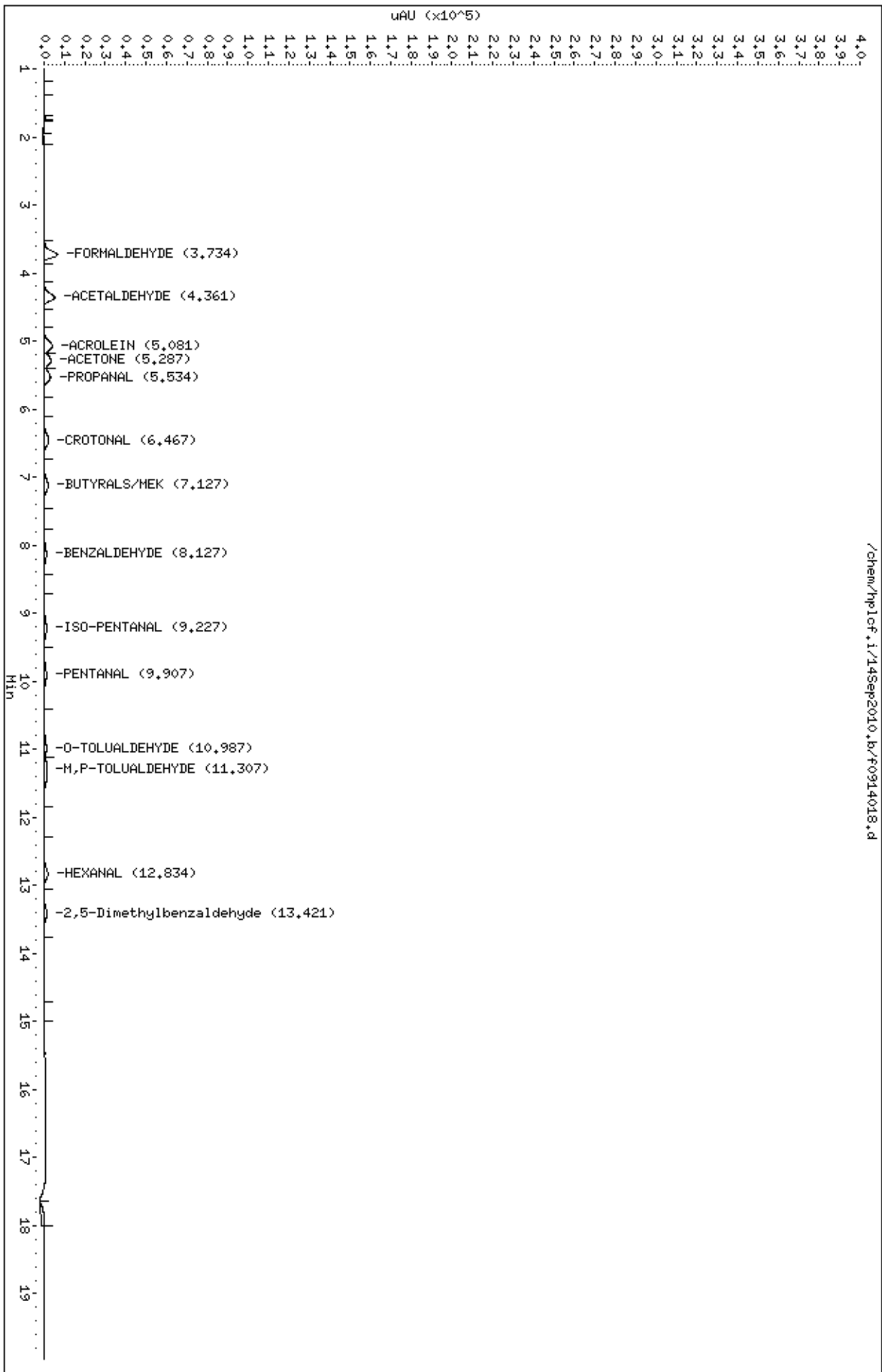
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds						AMOUNTS	
	RT	EXP RT	DLT RT	RT	RESPONSE	CAL - AMT	ON - COL
						(UG)	(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 ACROLEIN	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

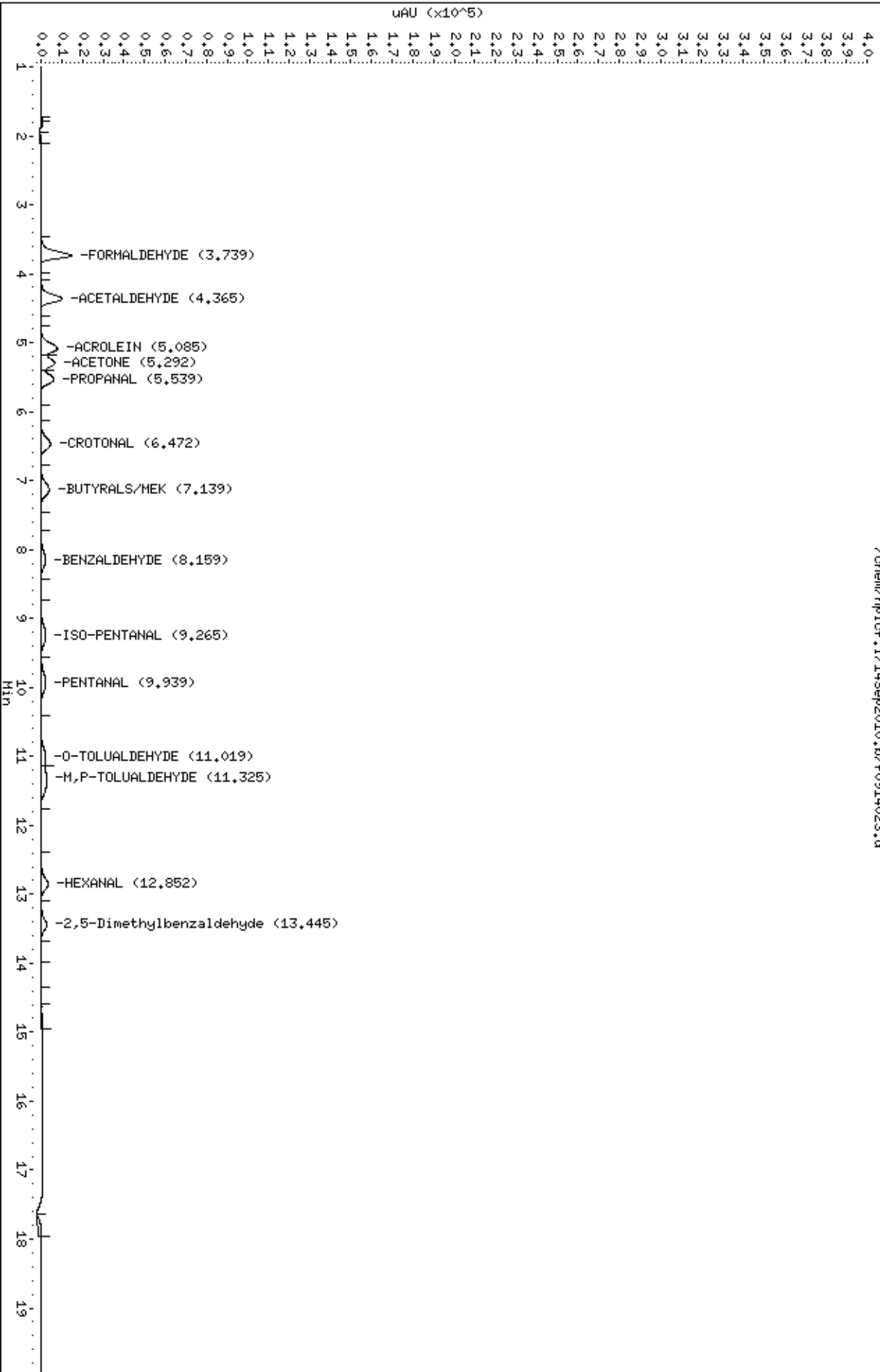
Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00

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



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds						AMOUNTS	
	RT	EXP RT	DLT RT	RT	RESPONSE	CAL - AMT	ON - COL
						(UG)	(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 ACROLEIN	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/hplcf.i/14Sep2010.b/f0914022.d

Date : 14-SEP-2010 21:17

Client ID: 5

Sample Info: #1838-34-1.0;5

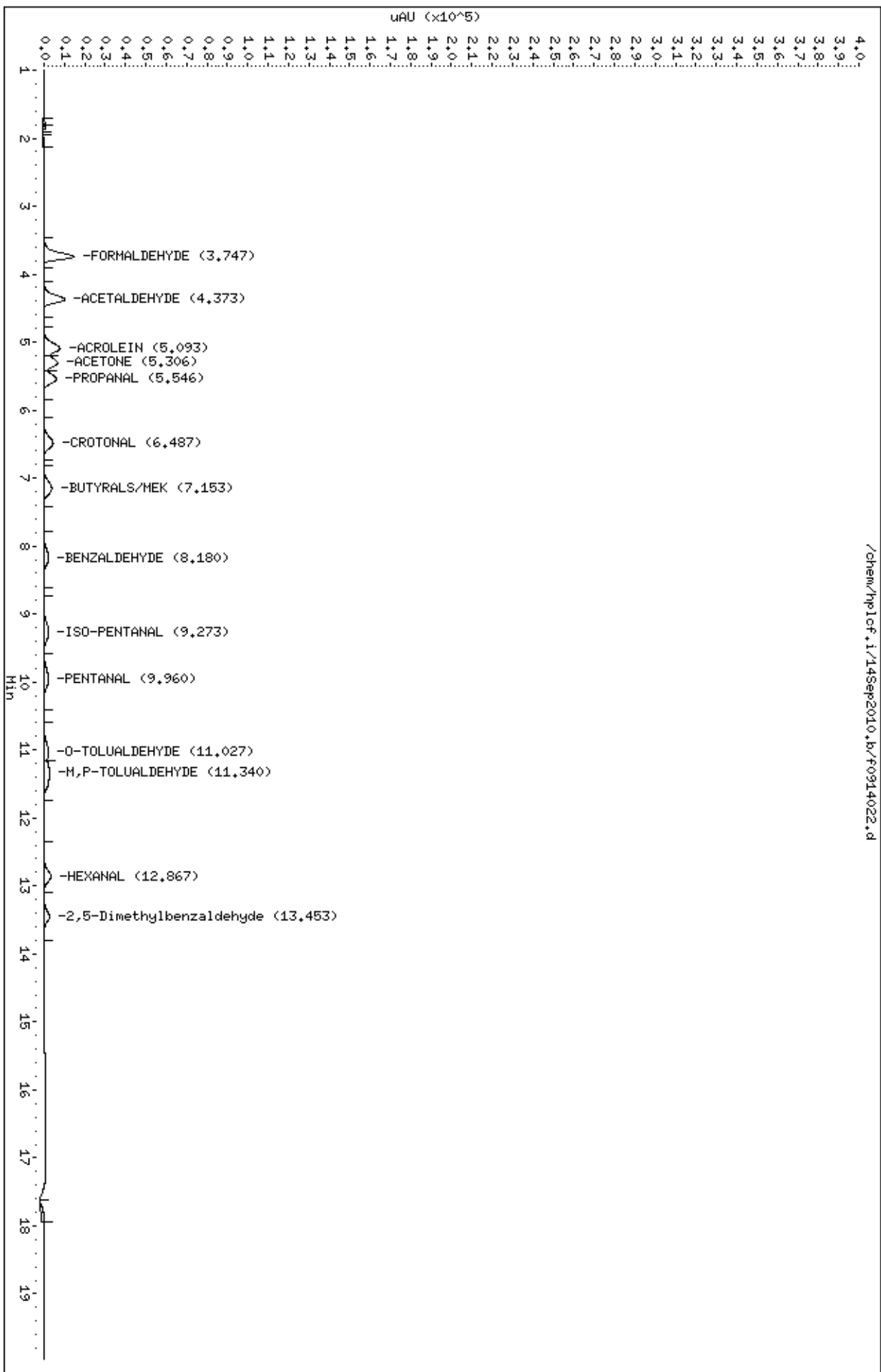
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds						AMOUNTS	
	RT	EXP RT	DLT RT	RT	RESPONSE	CAL - AMT	ON - COL
						(UG)	(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 ACROLEIN	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

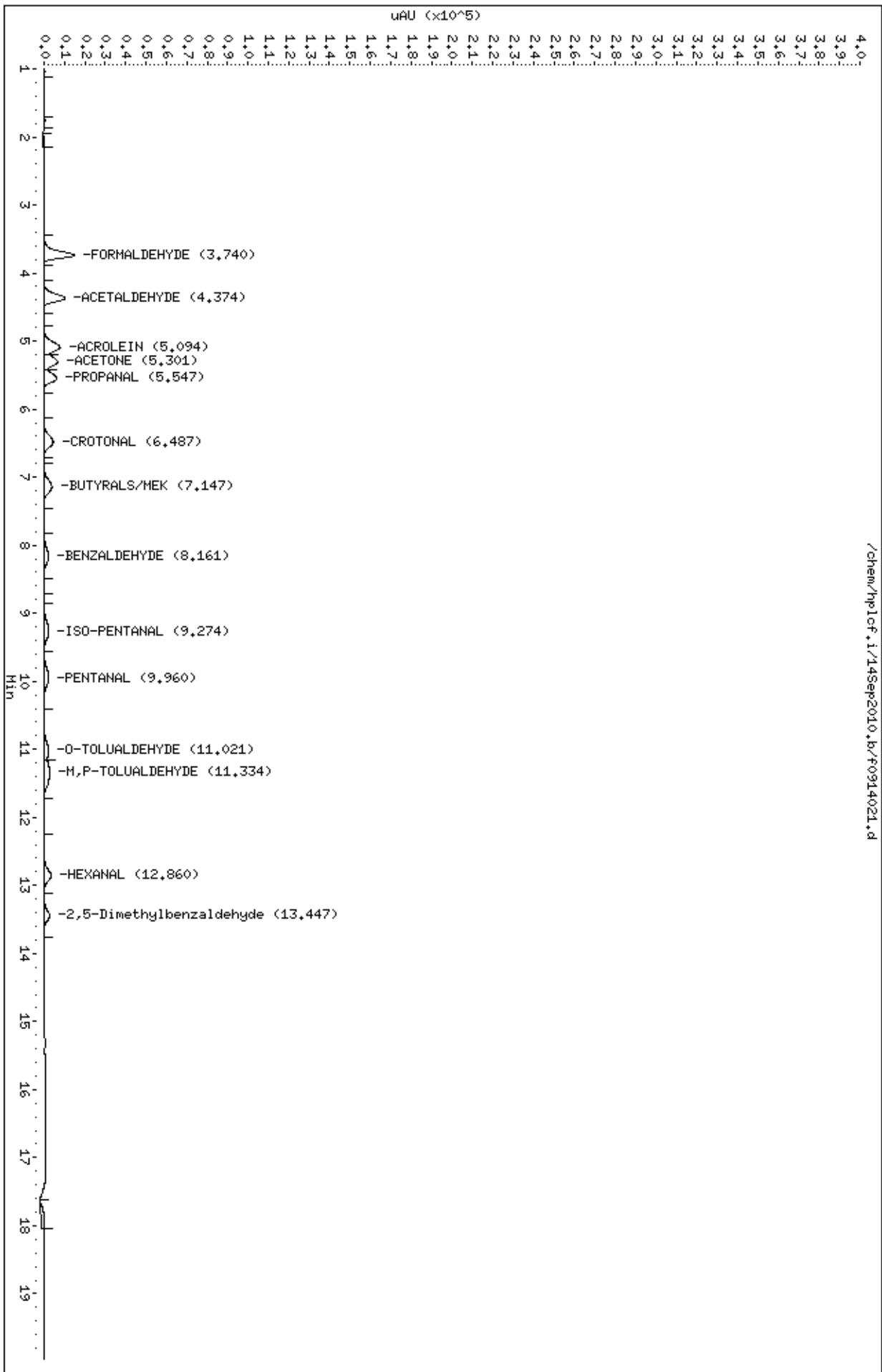
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds						AMOUNTS	
	RT	EXP RT	DLT RT	RT	RESPONSE	CAL - AMT	ON - COL
						(UG)	(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

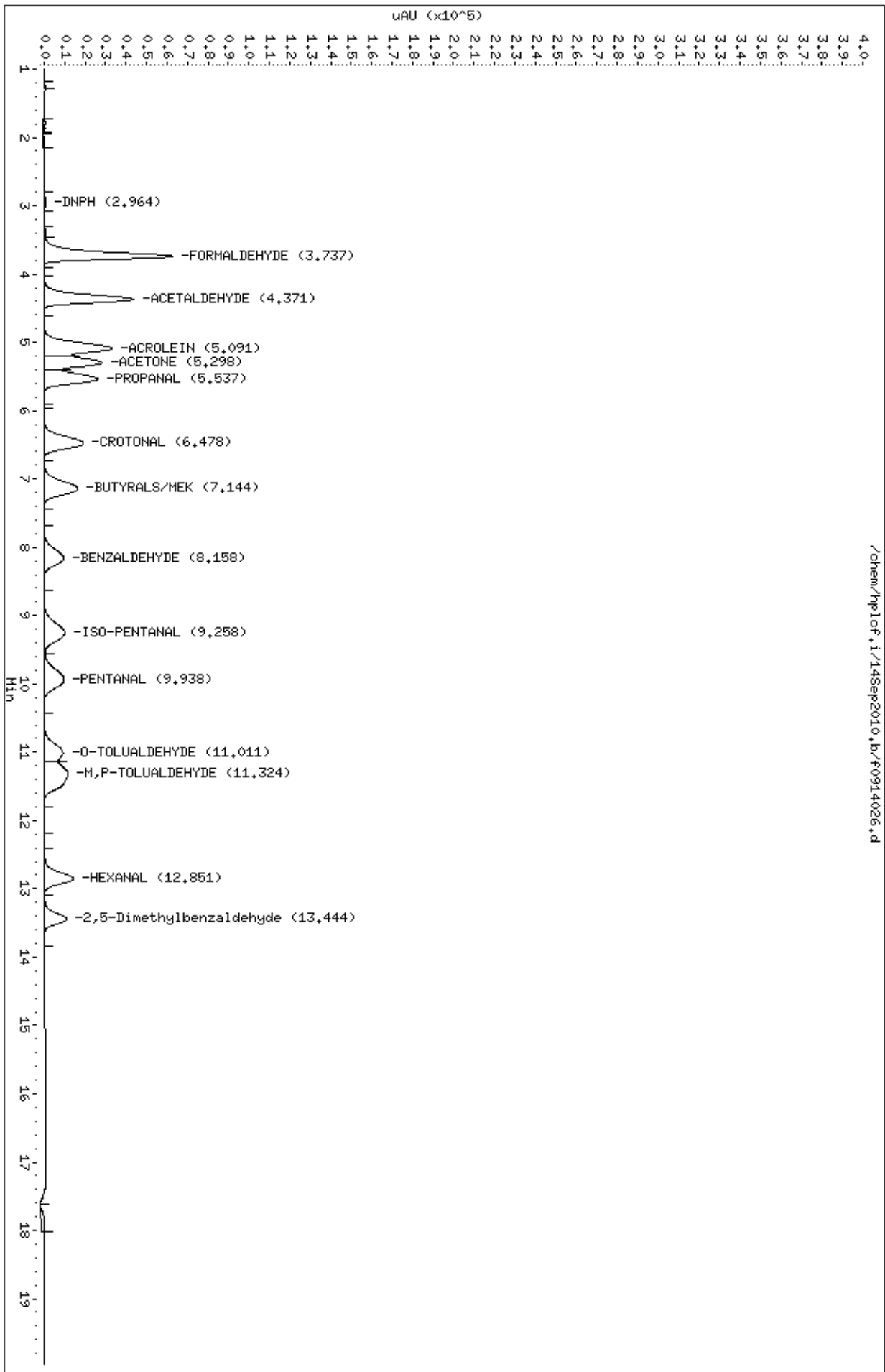
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds						AMOUNTS	
	RT	EXP RT	DLT RT	RT	RESPONSE	CAL - AMT	ON - COL
						(UG)	(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

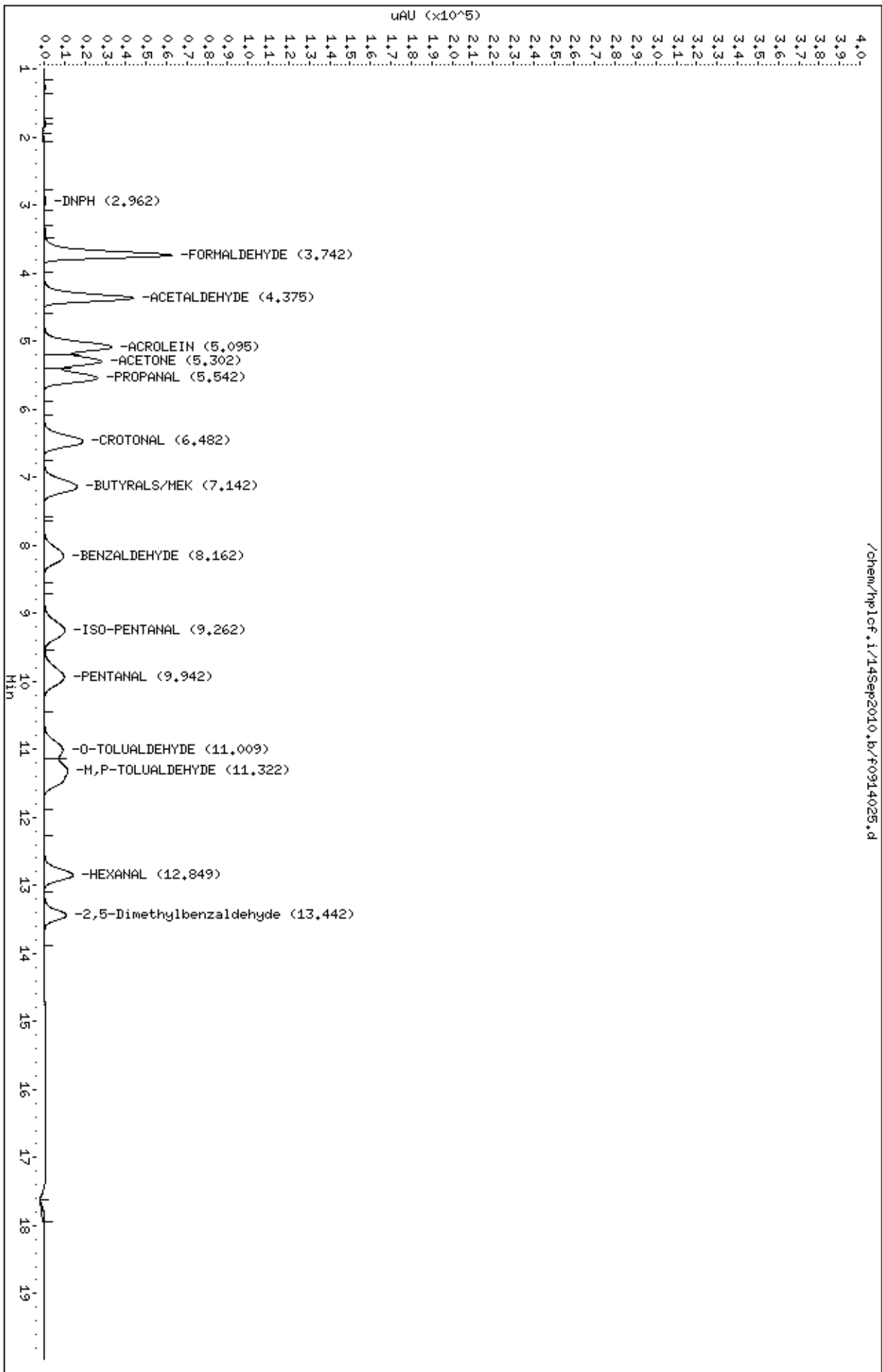
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/14Sep2010.b/f0914024.d

Lab Smp Id: 1838-32-4.0Client Smp ID: 6

Inj Date : 14-SEP-2010 21:59

Operator : CRLInst 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 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 21:59Cal File: f0914024.d

Als bottle: 1Calibration Sample, Level: 6

Dil Factor: 1.00000

Integrator: FalconCompound Sublist: all.sub

Target Version: 3.50

Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds						AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE		CAL - AMT	ON - COL
						(UG)	(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

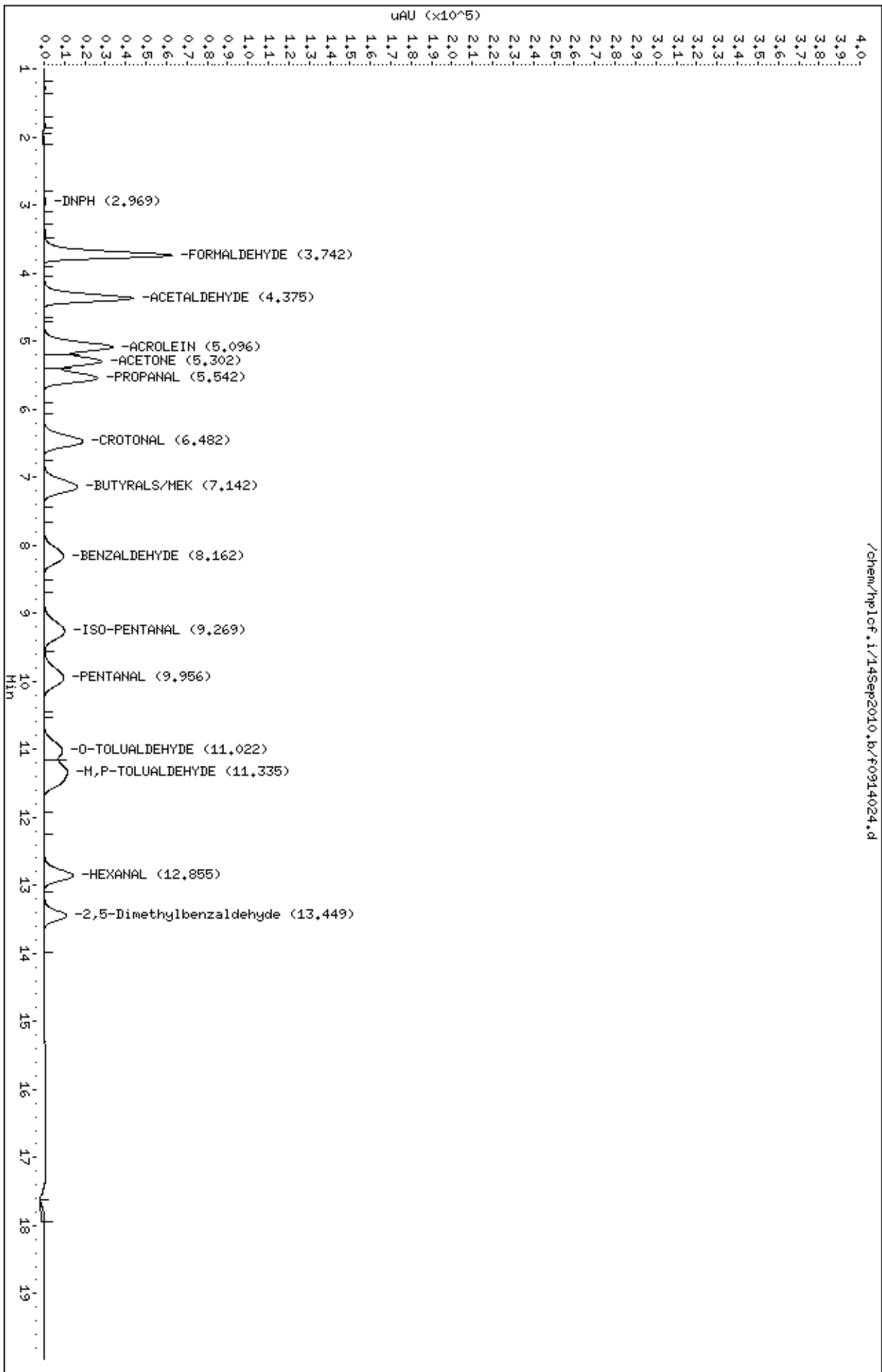
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds	AMOUNTS					
	RT			RESPONSE	CAL - AMT	ON - COL
	RT	EXP RT	DLT RT		(UG)	(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 AROLEIN	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

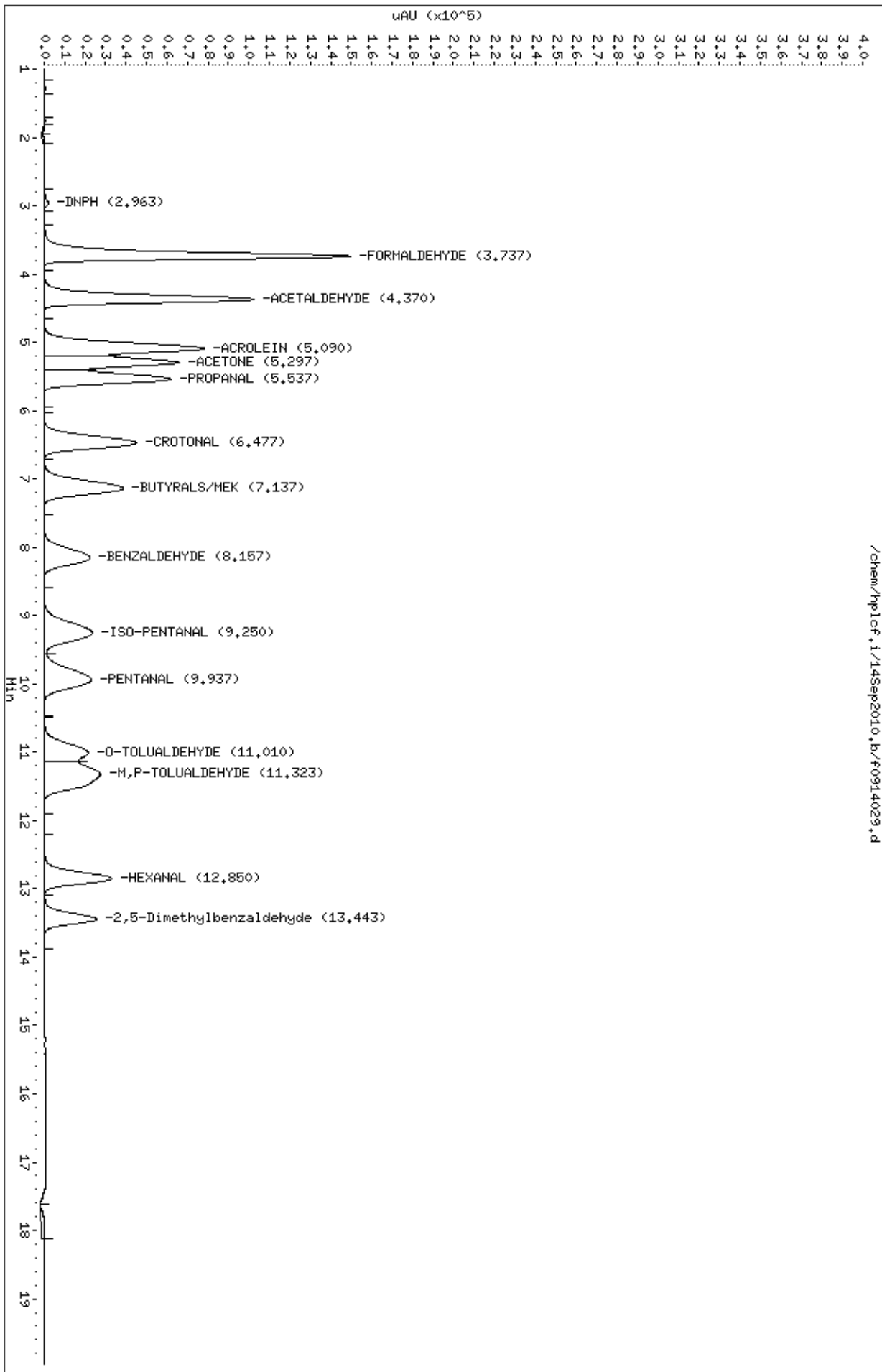
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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	AMOUNTS					
	RT			RESPONSE	CAL - AMT	ON - COL
	RT	EXP RT	DLT RT		(UG)	(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 AROLEIN	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: #1838-34-10.0:7

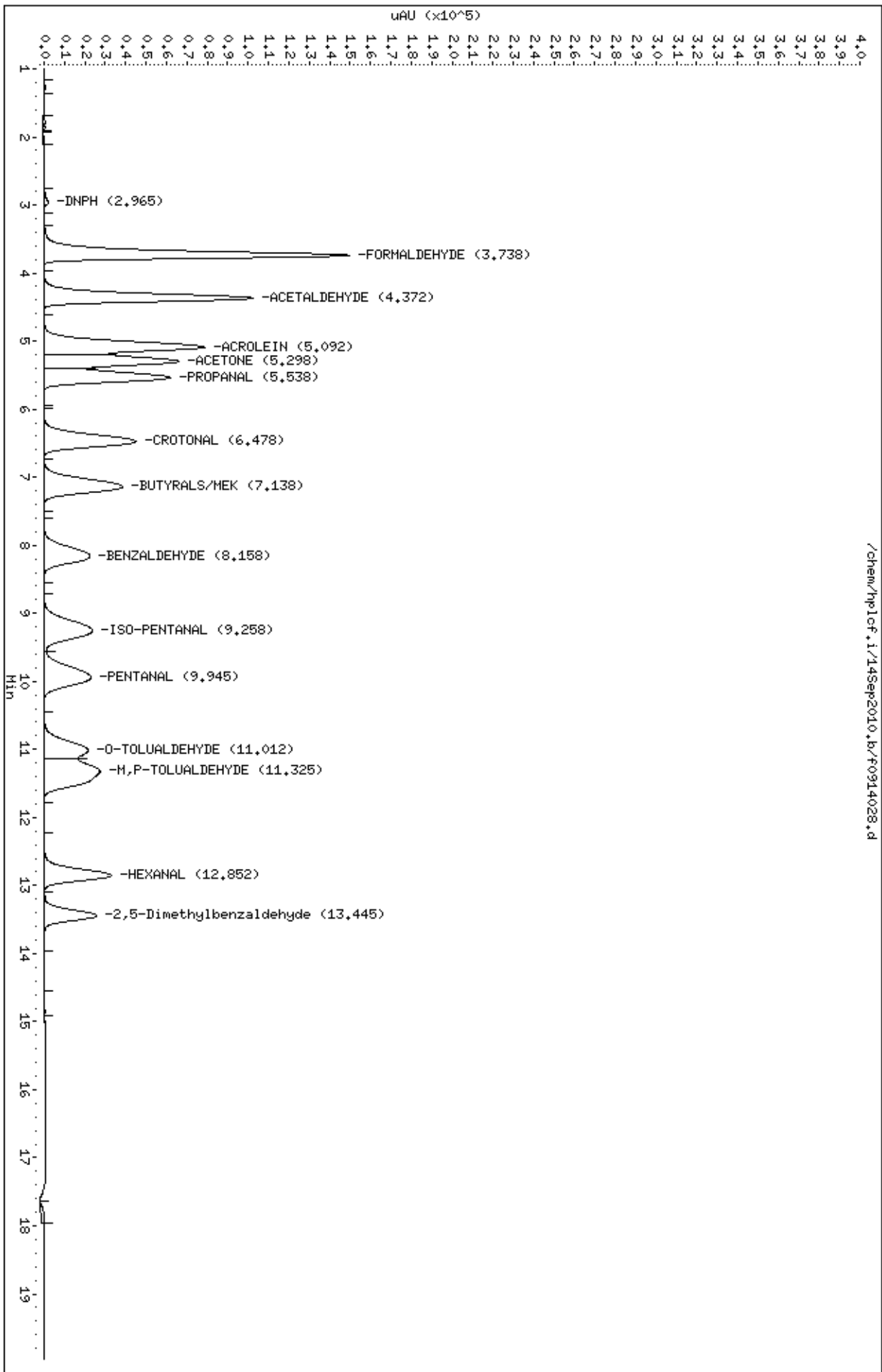
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/14Sep2010.b/f0914027.d

Lab Smp Id: 1838-34-10.0Client Smp ID: 7

Inj Date : 14-SEP-2010 23:02

Operator : CRLInst 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 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 23:02Cal File: f0914027.d

Als bottle: 1Calibration Sample, Level: 7

Dil Factor: 1.00000

Integrator: FalconCompound Sublist: all.sub

Target Version: 3.50

Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds	AMOUNTS					
	RT			RESPONSE	CAL - AMT	ON - COL
	RT	EXP RT	DLT RT		(UG)	(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 AROLEIN	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

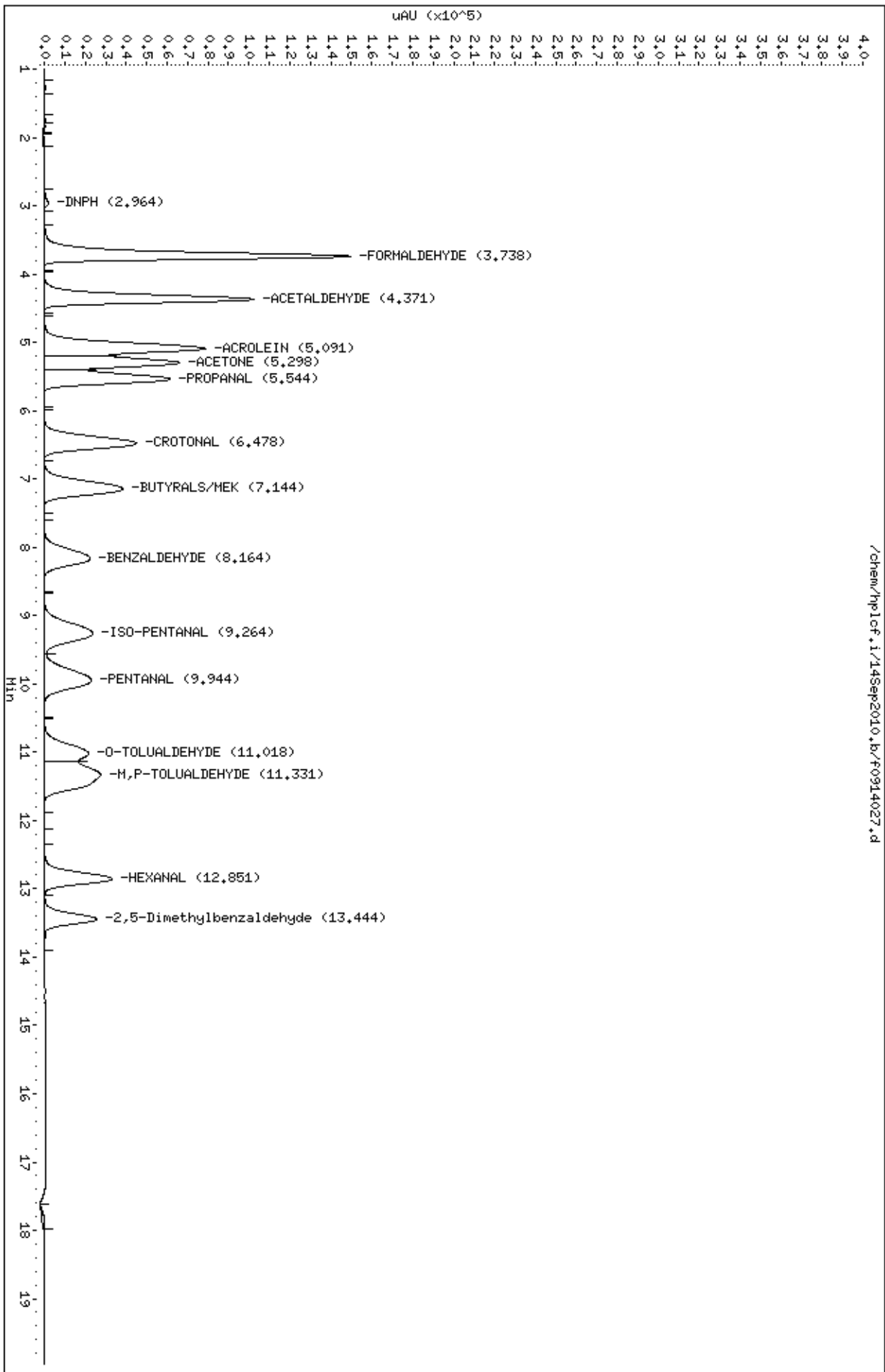
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

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

Lab Smp Id: 1838-34-20.0Client Smp ID: 8

Inj Date : 15-SEP-2010 00:46

Operator : CRLInst 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 cleafQuant Type: ESTD

Cal Date : 15-SEP-2010 00:46Cal File: f0914032.d

Als bottle: 1Calibration Sample, Level: 8

Dil Factor: 1.00000

Integrator: FalconCompound Sublist: all.sub

Target Version: 3.50

Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds						AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE		CAL - AMT	ON - COL
						(UG)	(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 AROLEIN	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/hp1cf.i/14Sep2010.b/f0914032.d

Date : 15-SEP-2010 00:46

Client ID: 8

Sample Info: J1838-34-20.0:8

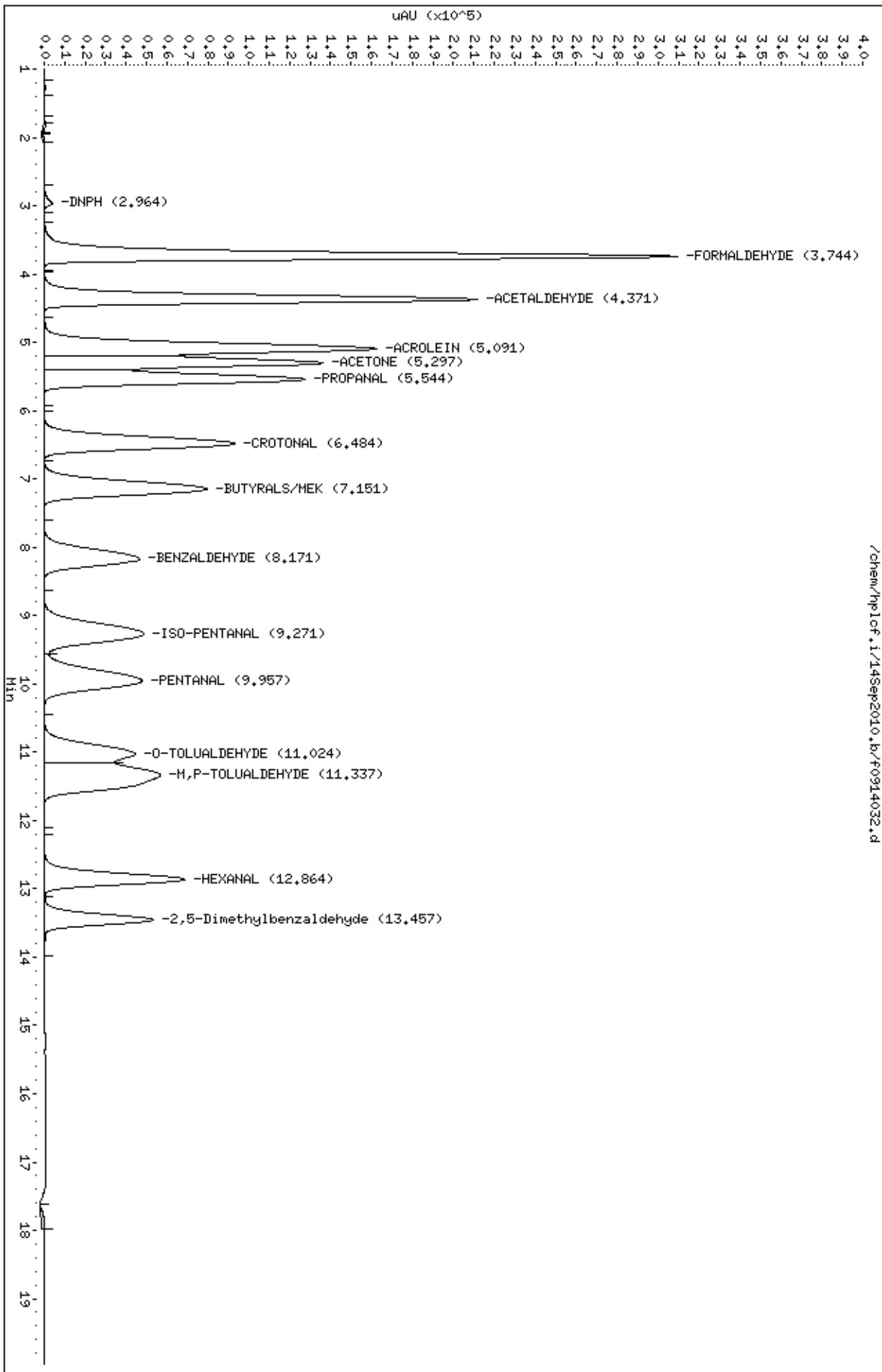
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/14Sep2010.b/f0914031.d

Lab Smp Id: 1838-34-20.0Client Smp ID: 8

Inj Date : 15-SEP-2010 00:25

Operator : CRLInst 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 cleafQuant Type: ESTD

Cal Date : 15-SEP-2010 00:25Cal File: f0914031.d

Als bottle: 1Calibration Sample, Level: 8

Dil Factor: 1.00000

Integrator: FalconCompound Sublist: all.sub

Target Version: 3.50

Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve 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	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

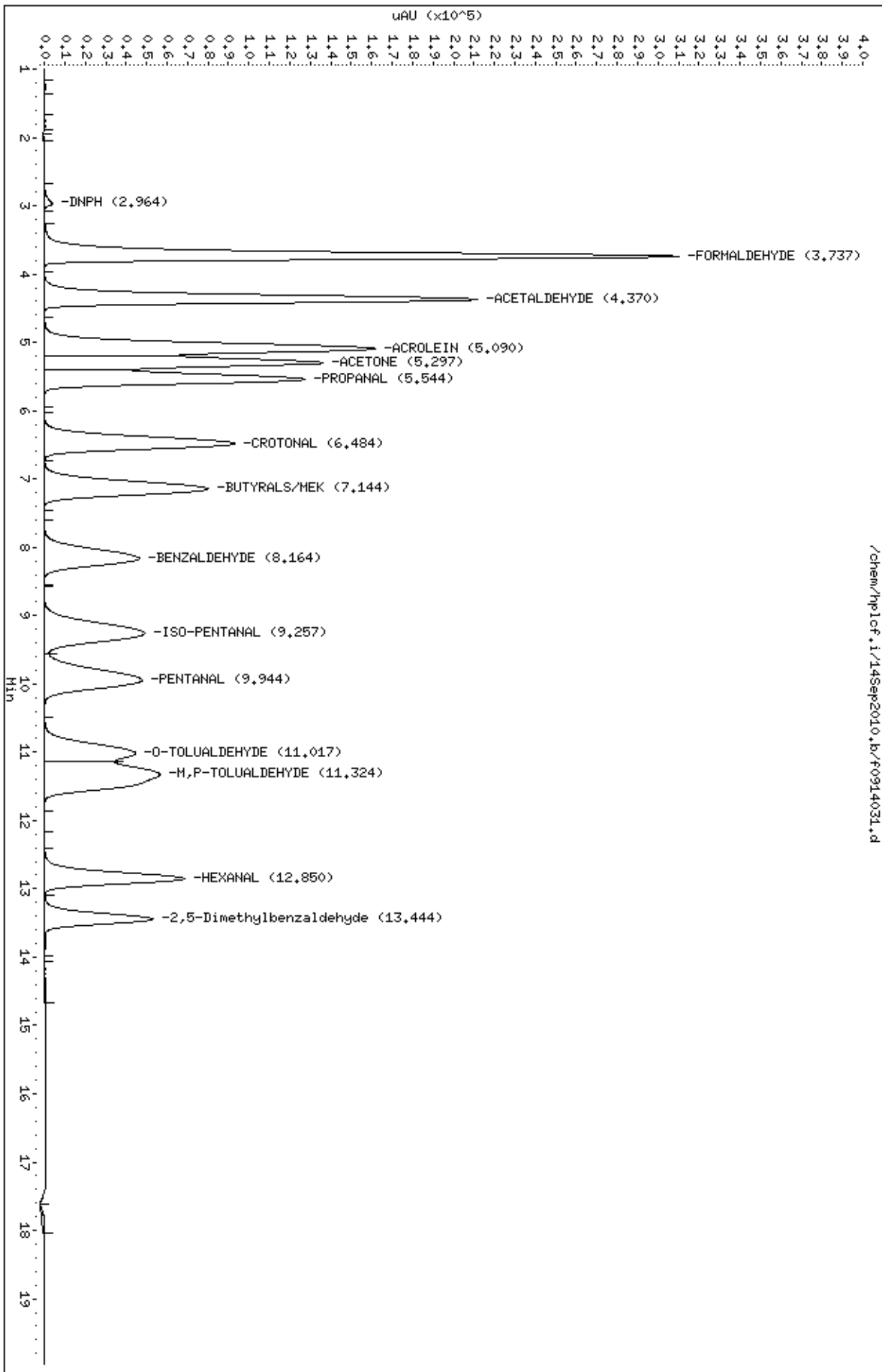
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-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	AMOUNTS					
	CAL - AMT			ON - COL		
	RT	EXP RT	DLT RT	RESPONSE	(UG)	(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

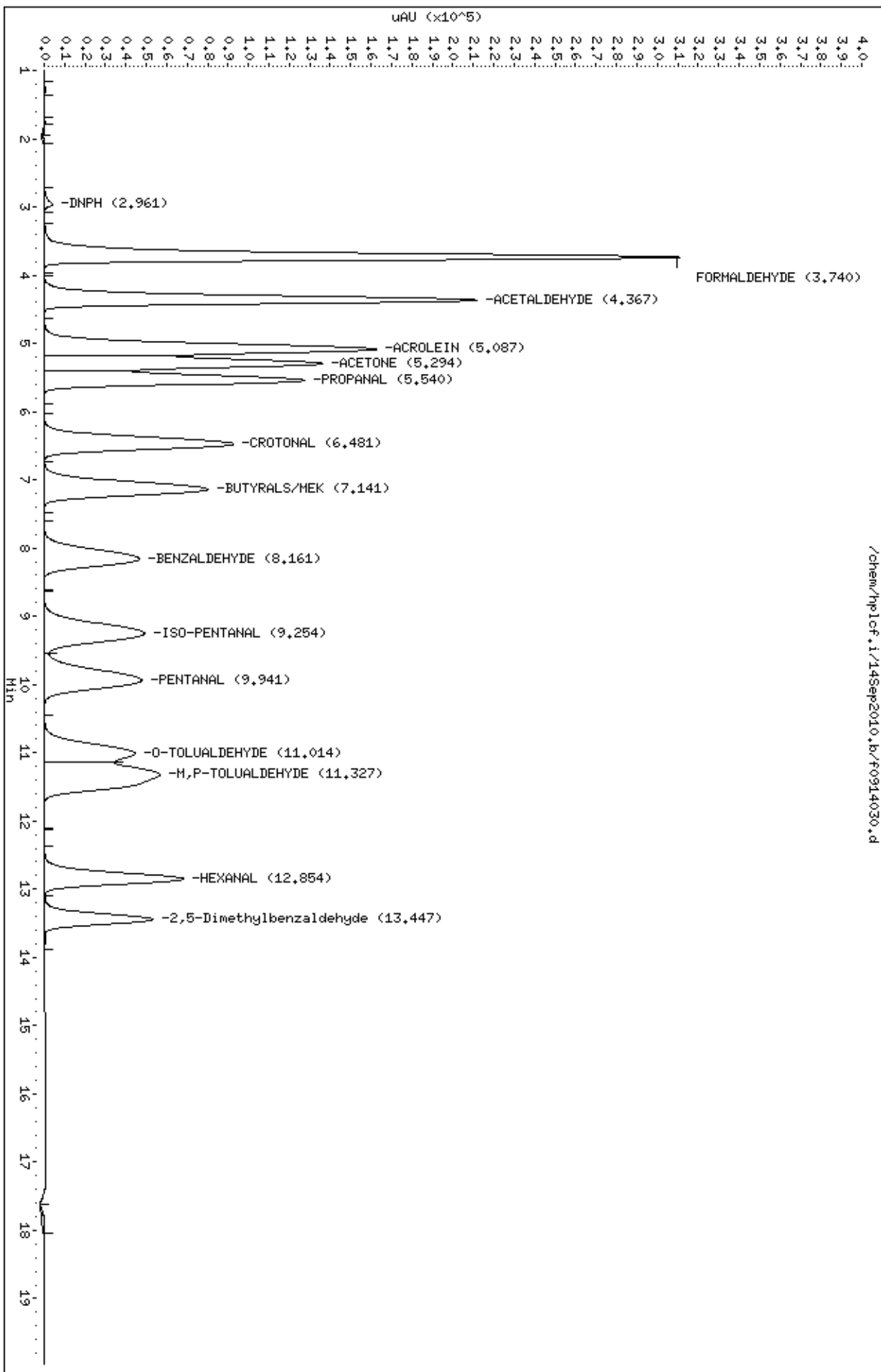
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

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

Lab Smp Id: 1838-34-50.0Client Smp ID: 9

Inj Date : 15-SEP-2010 01:49

Operator : CRLInst 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 cleafQuant Type: ESTD

Cal Date : 15-SEP-2010 01:49Cal File: f0914035.d

Als bottle: 1Calibration Sample, Level: 9

Dil Factor: 1.00000

Integrator: FalconCompound Sublist: all.sub

Target Version: 3.50

Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL - AMT	ON - COL
					(UG)	(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 ACROLEIN	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: J1838-34-50.0:9

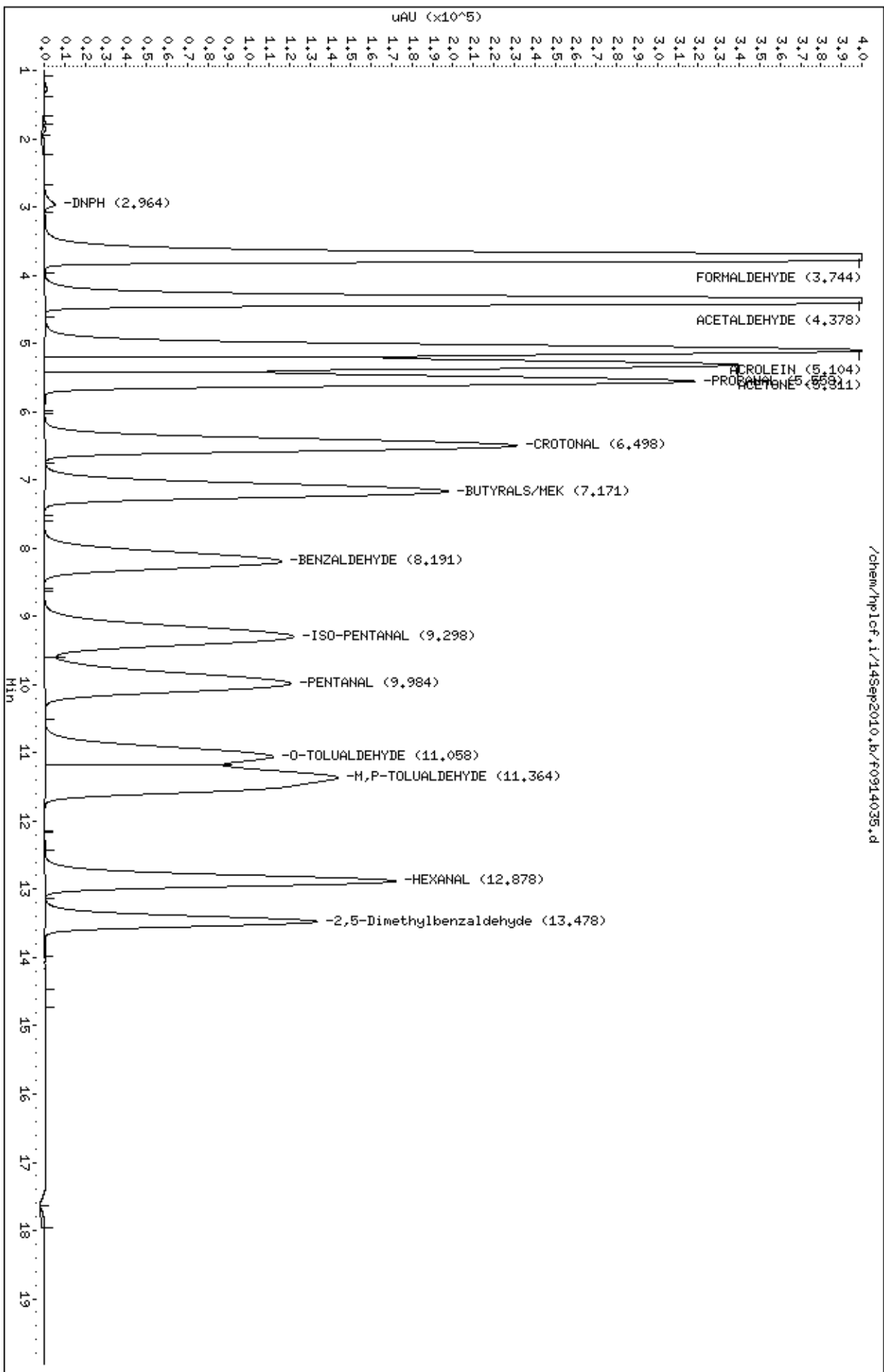
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

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

Lab Smp Id: 1838-34-50.0Client Smp ID: 9

Inj Date : 15-SEP-2010 01:28

Operator : CRLInst 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 cleafQuant Type: ESTD

Cal Date : 15-SEP-2010 01:28Cal File: f0914034.d

Als bottle: 1Calibration Sample, Level: 9

Dil Factor: 1.00000

Integrator: FalconCompound Sublist: all.sub

Target Version: 3.50

Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds						AMOUNTS	
	RT	EXP RT	DLT RT	RT	RESPONSE	CAL - AMT	ON - COL
						(UG)	(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 ACROLEIN	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

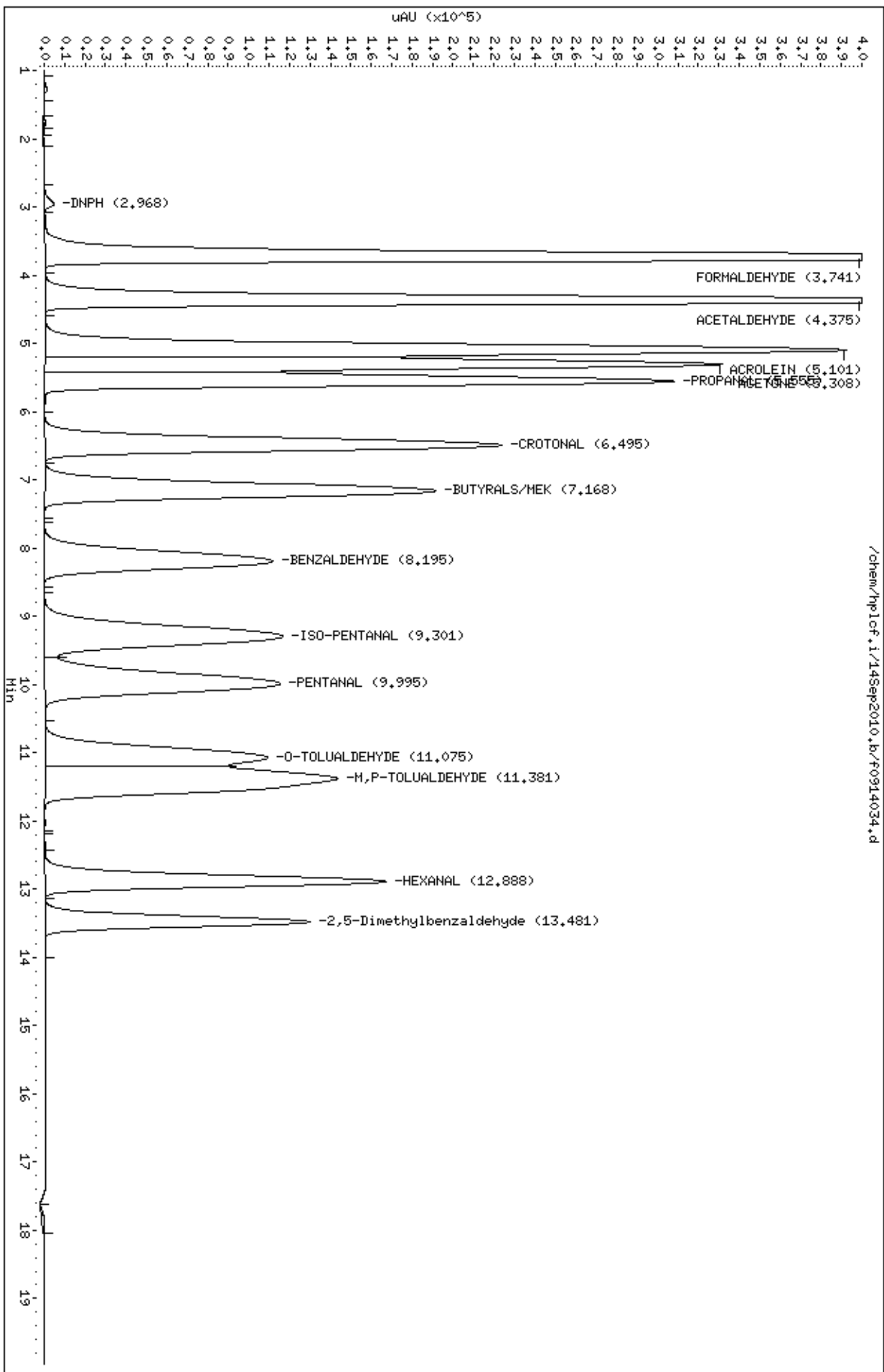
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

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

Lab Smp Id: 1838-34-50.0Client Smp ID: 9

Inj Date : 15-SEP-2010 01:07

Operator : CRLInst 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 cleafQuant Type: ESTD

Cal Date : 15-SEP-2010 01:07Cal File: f0914033.d

Als bottle: 1Calibration Sample, Level: 9

Dil Factor: 1.00000

Integrator: FalconCompound Sublist: all.sub

Target Version: 3.50

Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL - AMT	ON - COL
					(UG)	(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 ACROLEIN	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

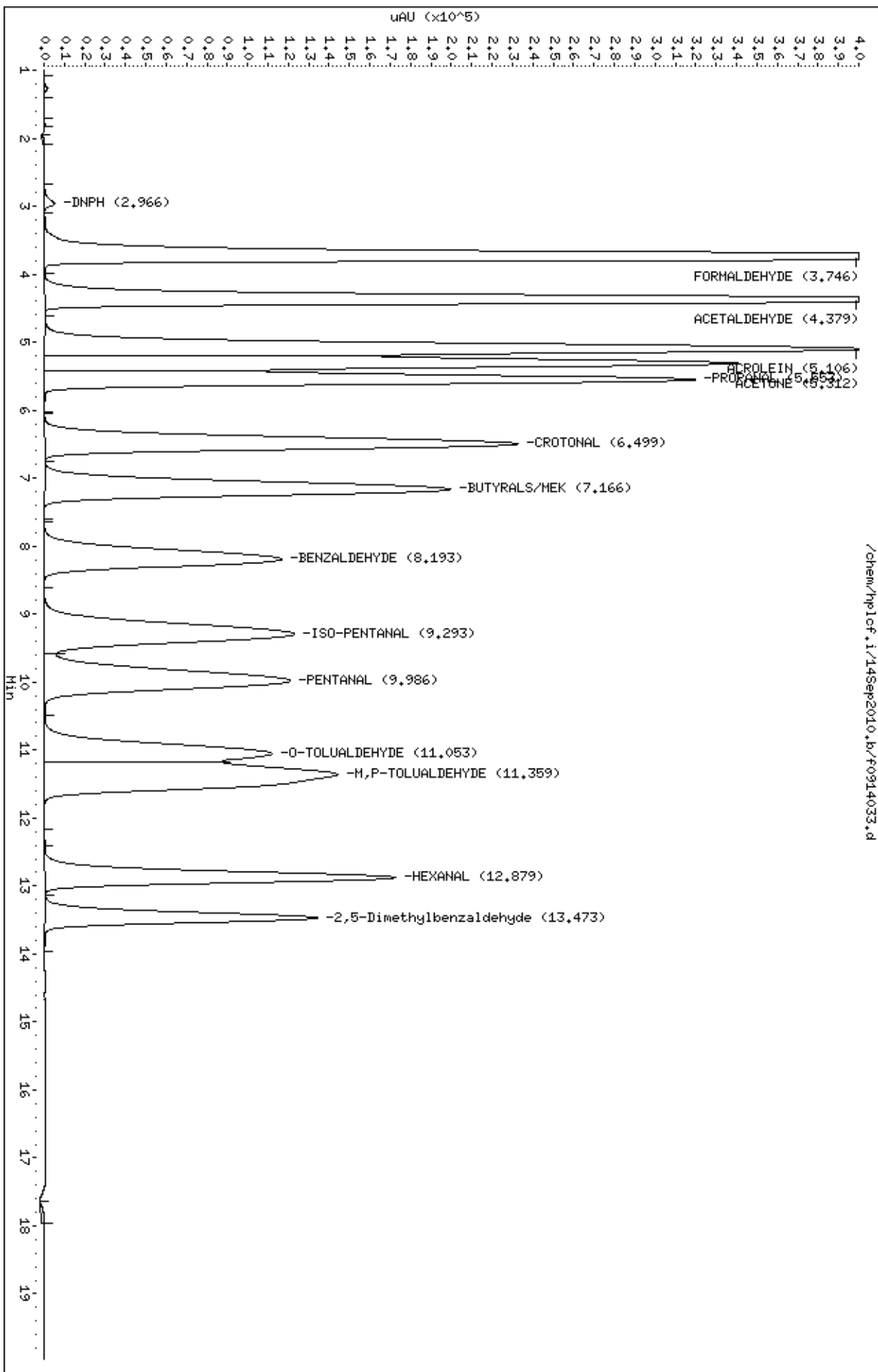
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/14Sep2010.b/f0914038.d

Lab Smp Id: 1838-34-75.0Client Smp ID: 10

Inj Date : 15-SEP-2010 02:52

Operator : CRLInst 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 cleafQuant Type: ESTD

Cal Date : 15-SEP-2010 02:52Cal File: f0914038.d

Als bottle: 1Calibration Sample, Level: 10

Dil Factor: 1.00000

Integrator: FalconCompound Sublist: all.sub

Target Version: 3.50

Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL - AMT	ON - COL
					(UG)	(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 AROLEIN	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

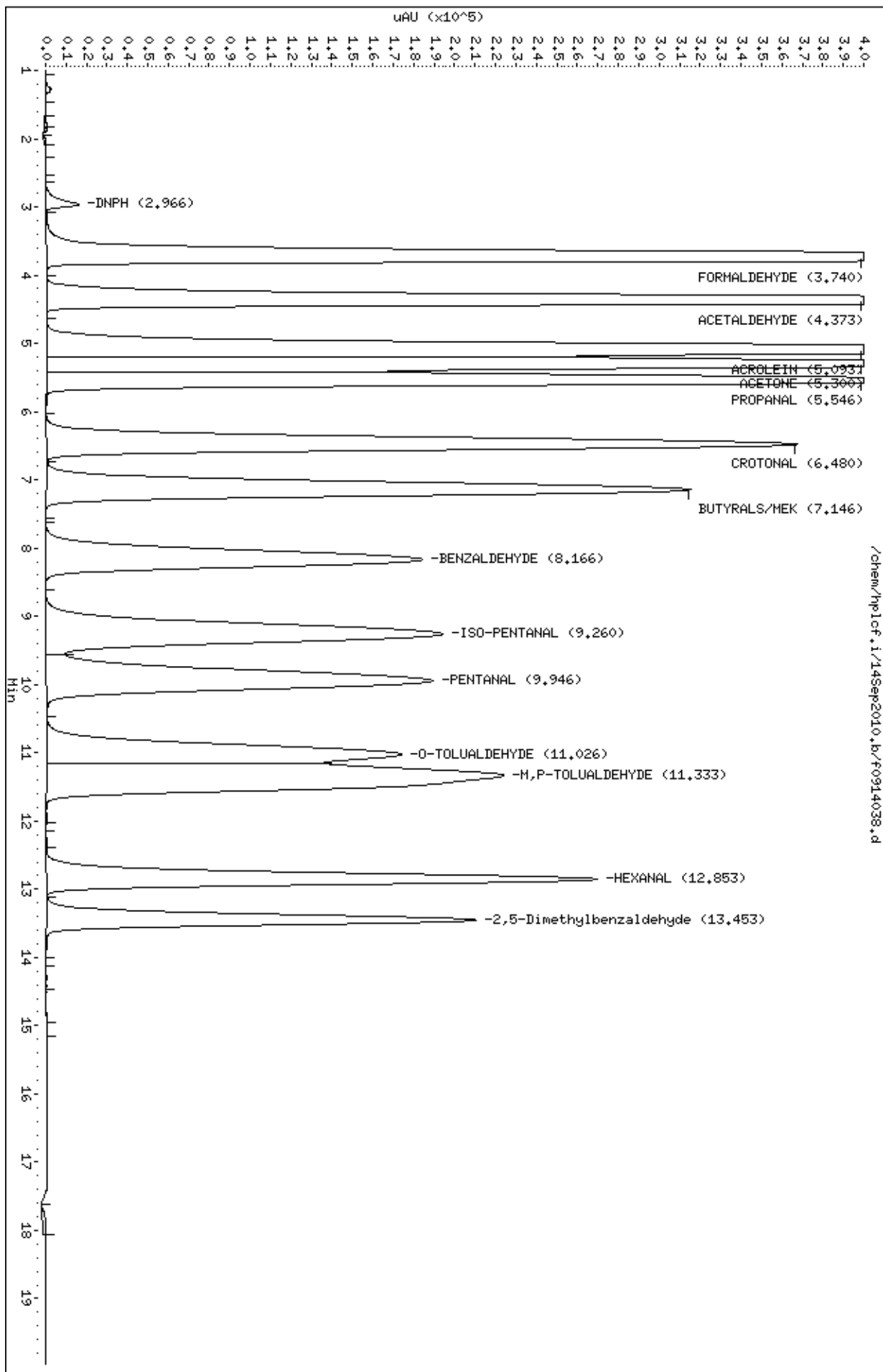
Purge Volume: 5.0

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

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds	AMOUNTS					
	RT			RESPONSE	CAL - AMT	ON - COL
	RT	EXP RT	DLT RT		(UG)	(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

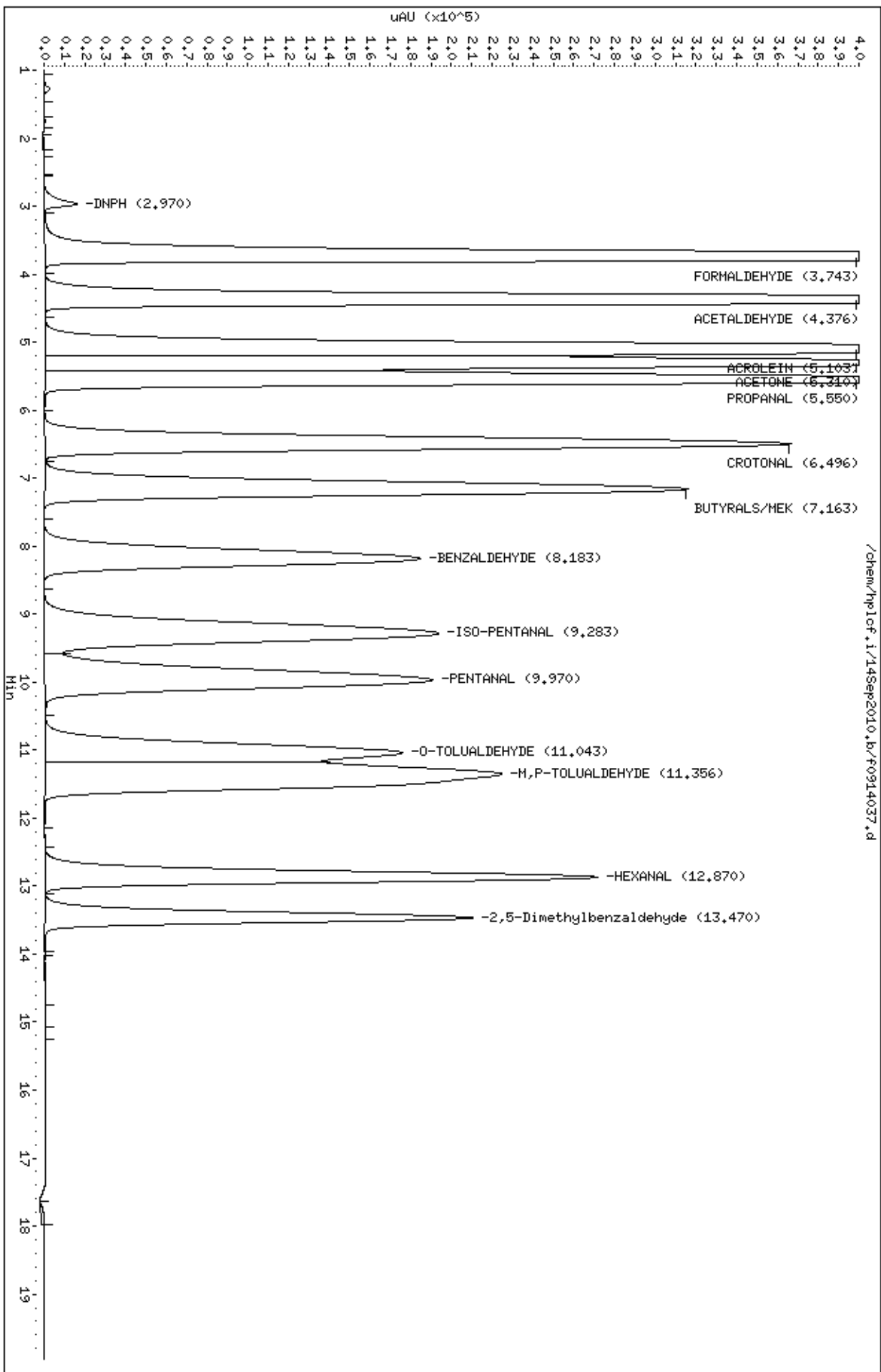
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hp1cf.i

Operator: CRL

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/14Sep2010.b/f0914036.d

Lab Smp Id: 1838-34-75.0Client Smp ID: 10

Inj Date : 15-SEP-2010 02:10

Operator : CRLInst 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 cleafQuant Type: ESTD

Cal Date : 15-SEP-2010 02:10Cal File: f0914036.d

Als bottle: 1Calibration Sample, Level: 10

Dil Factor: 1.00000

Integrator: FalconCompound Sublist: all.sub

Target Version: 3.50

Processing Host: eeyore

Concentration Formula: Amt * DF * Vs/Vo * CpndVariable

Vs5.00000Sample Final Volume (mL)

Vo5.00000Curve based Volume (mL)

Cpnd Variable

Local Compound Variable

Compounds					AMOUNTS	
	RT	EXP RT	DLT RT	RESPONSE	CAL - AMT	ON - COL
					(UG)	(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 ACROLEIN	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

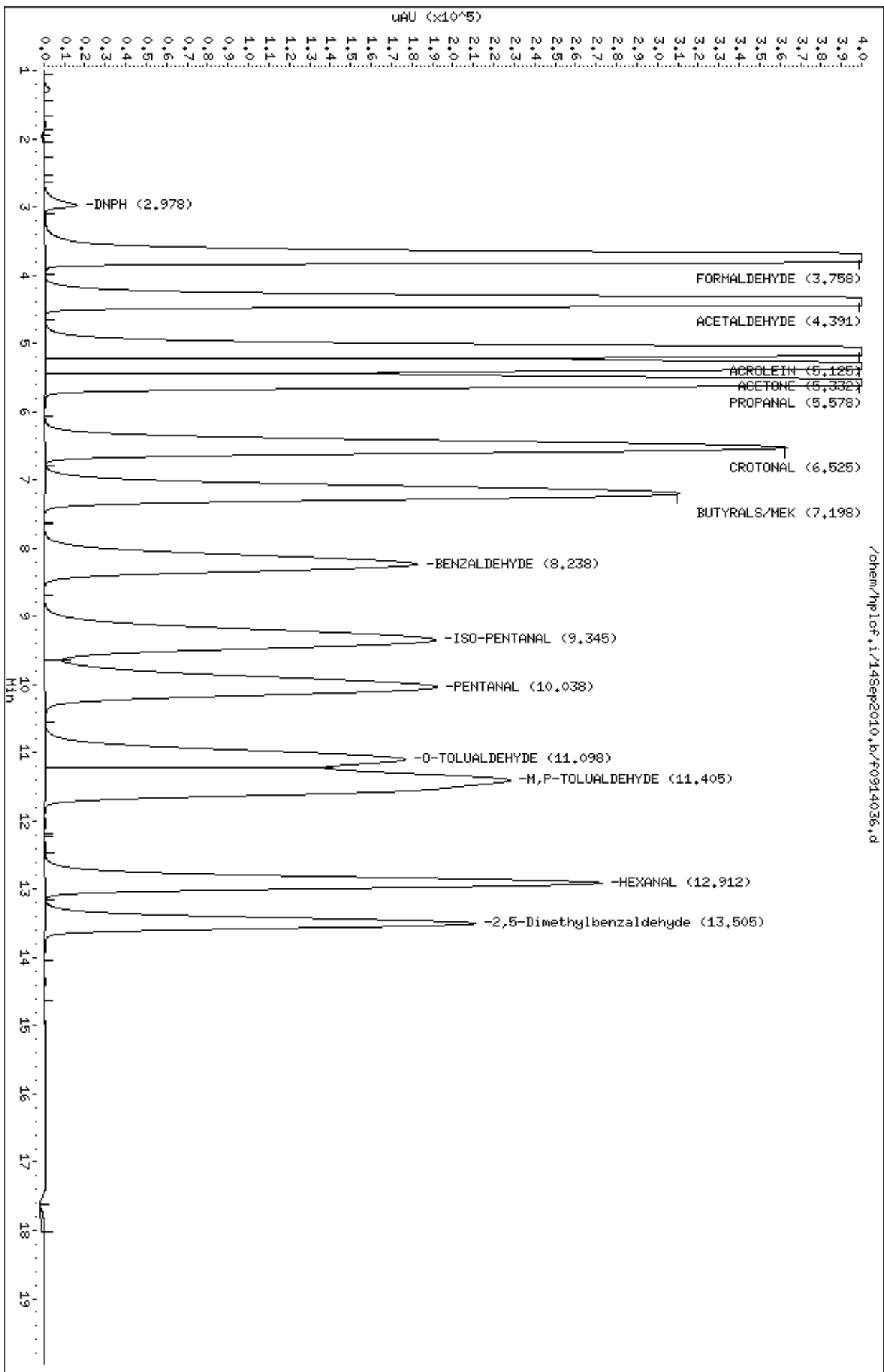
Purge Volume: 5.0

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: 23-NOV-2010 08:01

Lab File ID: f1123002.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/23Nov2010.b/F10D0914.m

COMPOUND	RRF / AMOUNT	RF10	MIN RRF	%D / %DRIFT	MAX %D / %DRIFT	CURVE TYPE
1 DNPH	++++	526	0.100	++++	10.00000	Averaged<-
2 FORMALDEHYDE	109474	113650	0.010	-3.81400	10.00000	Averaged
3 ACETALDEHYDE	85159	85114	0.010	0.05226	10.00000	Averaged
4 ACRROLEIN	81132	86449	0.010	-6.55451	10.00000	Averaged
5 ACETONE	62324	61577	0.010	1.19905	10.00000	Averaged
6 PROPANAL	62684	56890	0.010	9.24348	10.00000	Averaged
7 CROTONAL	56944	55890	0.010	1.85035	10.00000	Averaged
9 BUTYRALS/MEK	54133	52013	0.010	3.91657	10.00000	Averaged
10 BENZALDEHYDE	38259	37230	0.010	2.68884	10.00000	Averaged
11 ISO-PENTANAL	45456	45909	0.010	-0.99648	10.00000	Averaged
12 PENTANAL	44700	42942	0.010	3.93313	10.00000	Averaged
14 O-TOLUALDEHYDE	36466	36574	0.010	-0.29634	10.00000	Averaged
15 M,P-TOLUALDEHYDE	30634	28921	0.010	5.59255	10.00000	Averaged
16 HEXANAL	40334	39534	0.010	1.98448	10.00000	Averaged
17 2,5-Dimethylbenzaldehyde	29495	28228	0.010	4.29589	10.00000	Averaged

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123002.d

Lab Smp Id: 1838-41-10Client Smp ID: CCV

Inj Date : 23-NOV-2010 08:01

Operator : FMInst ID: hplcf.i

Smp Info : ;1838-41-10;CCV

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 23-Nov-2010 09:21 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1Continuing Calibration Sample

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	AMOUNTS					
	RT	EXP RT	DLT RT	RESPONSE	CAL-AMT (ng)	ON-COL (ng)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.598	3.598	0.000	1136496	10.0000	10.4
3 ACETALDEHYDE	4.165	4.165	0.000	851141	10.0000	9.99
4 ACRROLEIN	4.791	4.791	0.000	864493	10.0000	10.6
5 ACETONE	4.991	4.991	0.000	615769	10.0000	9.88
6 PROPANAL	5.198	5.198	0.000	568899	10.0000	9.08
7 CROTONAL	5.991	5.991	0.000	558904	10.0000	9.81
9 BUTYRALS/MEK	6.565	6.565	0.000	520131	10.0000	9.61
10 BENZALDEHYDE	7.405	7.405	0.000	372301	10.0000	9.73
11 ISO-PENTANAL	8.338	8.338	0.000	459090	10.0000	10.1
12 PENTANAL	8.898	8.898	0.000	429421	10.0000	9.61
14 O-TOLUALDEHYDE	9.985	9.985	0.000	365742	10.0000	10.0
15 M,P-TOLUALDEHYDE	10.338	10.338	0.000	578411	20.0000	18.9
16 HEXANAL	12.091	12.091	0.000	395336	10.0000	9.80
17 2,5-Dimethylbenzaldehyde	12.745	12.745	0.000	282278	10.0000	9.57

Data File: /chem/hplcf.i/23Nov2010.b/f1123002.d

Date : 23-NOV-2010 08:01

Client ID: CCV

Sample Info: f1838-41-10;CCV

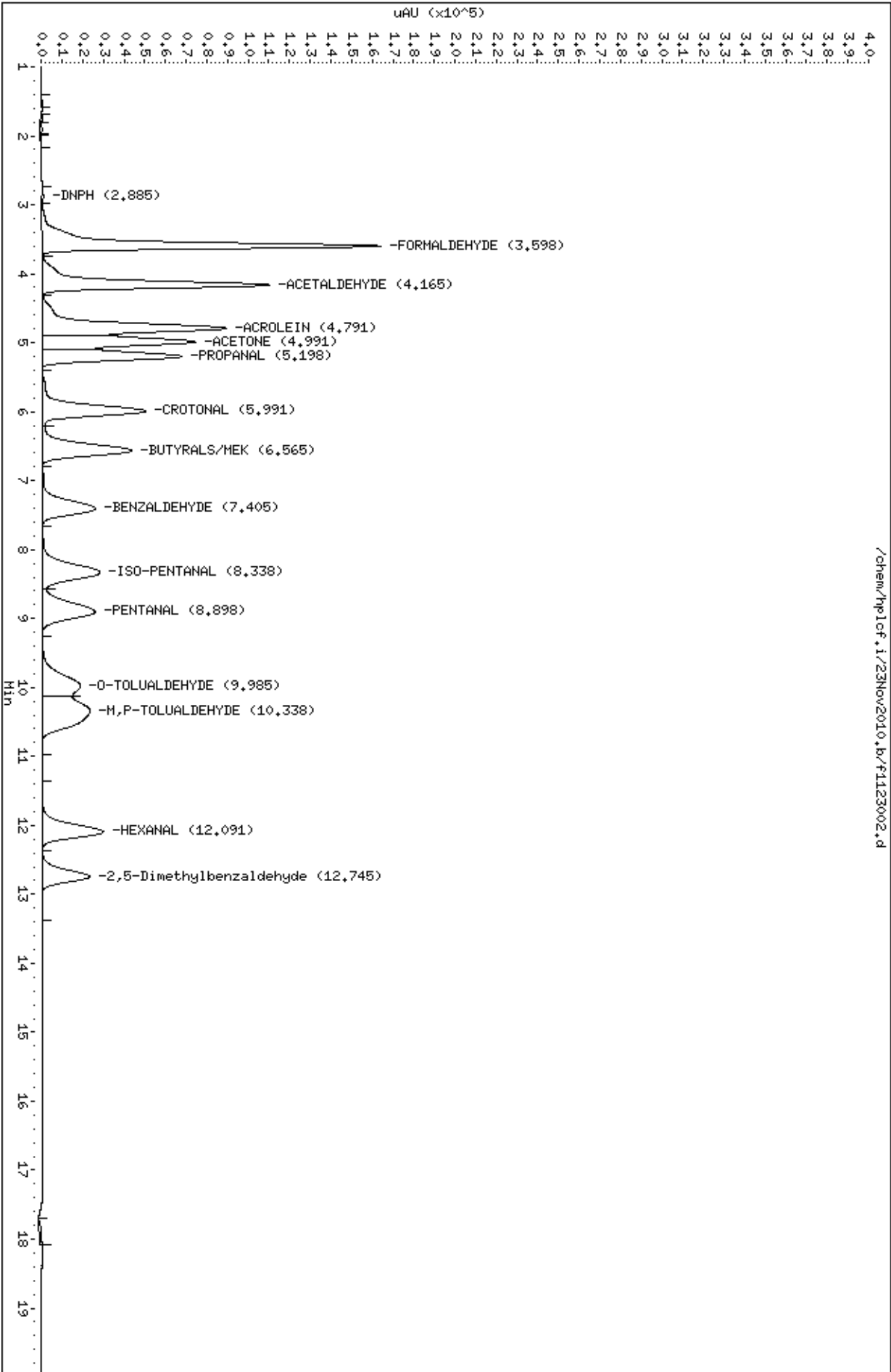
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: FM

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123003.d

Lab Smp Id: CCBClient Smp ID: LAB BLANK

Inj Date : 23-NOV-2010 08:22

Operator : FMInst ID: hplcf.i

Smp Info : ;CCB;LAB BLANK

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 23-Nov-2010 09:21 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
	RT	EXP RT	DLT RT	RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====		=====	=====	=====
1 DNPH					Compound Not Detected.		
2 FORMALDEHYDE					Compound Not Detected.		
3 ACETALDEHYDE					Compound Not Detected.		
4 ACRYLEIN					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	CONCENTRATIONS					
	RT	EXP RT	DLT	RT	RESPONSE	ON-COLUMN
						FINAL
=====	==	=====	=====		=====	(ng) (ug)
17 2,5-Dimethylbenzaldehyde						=====
					Compound Not Detected.	

Data File: /chem/hplcf.i/23Nov2010.b/f1123003.d

Page 1

Date : 23-NOV-2010 08:22

Client ID: LAB BLANK

Instrument: hplcf.i

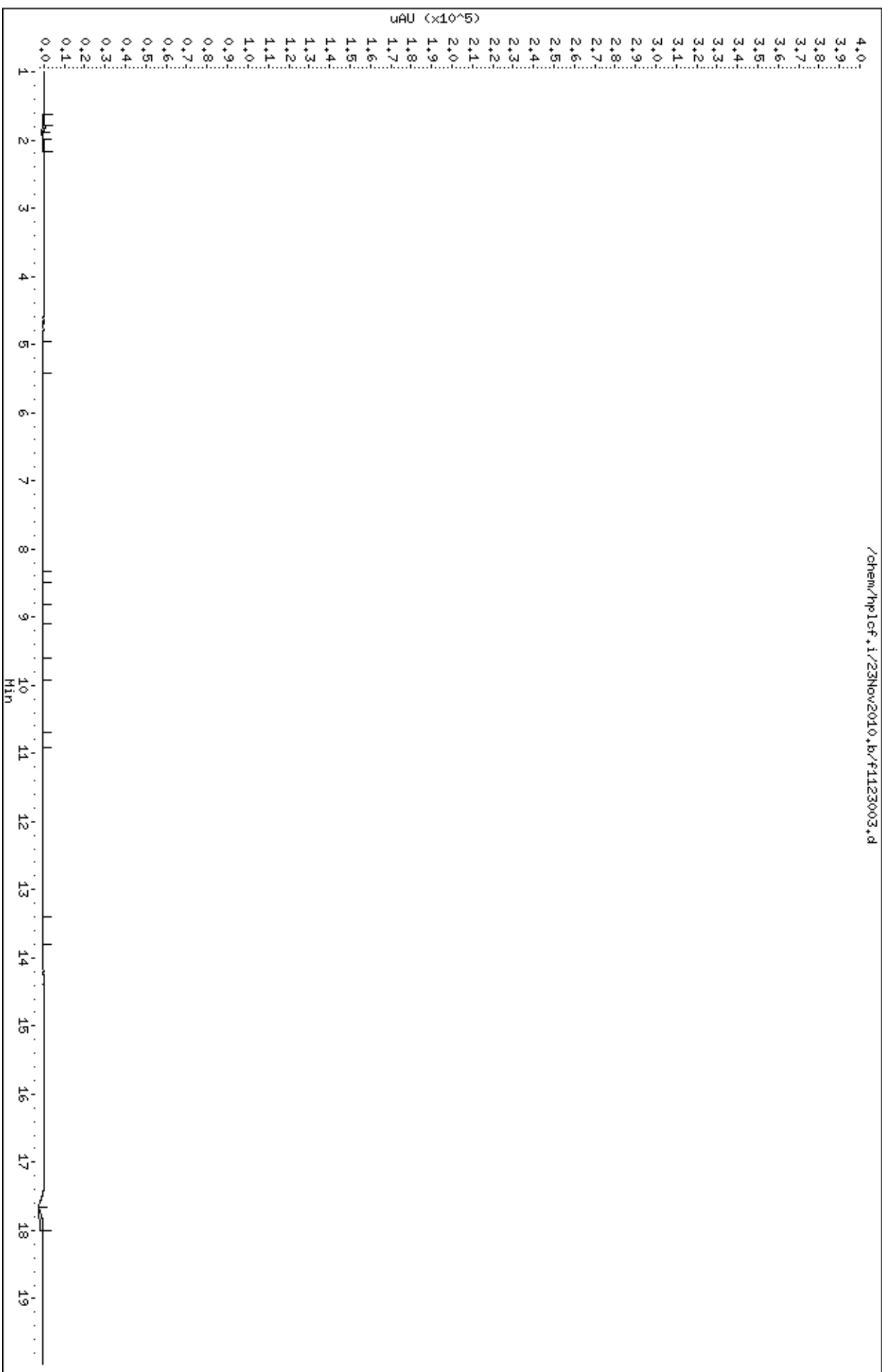
Sample Info: JCCB;LAB BLANK

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



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: 23Nov2010
Client Smp ID: CCV
Sample Point:
Date Received:
Quant Type: ESTD
Level: MED
Operator: FM

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/KG) ug	Q
119-26-6-----	DNPH	0.00	
50-00-0-----	FORMALDEHYDE	10.3	
75-07-0-----	ACETALDEHYDE	9.23	
107-02-8-----	ACROLEIN	10.1	
67-64-1-----	ACETONE	9.79	
123-38-6-----	PROPANAL	9.24	
123-73-9-----	CROTONAL	9.82	
9999-9999-007--	BUTYRALS/MEK	9.60	
100-52-7-----	BENZALDEHYDE	9.60	
590-86-3-----	ISO-PENTANAL	9.97	
110-62-3-----	PENTANAL	9.70	
529-20-4-----	O-TOLUALDEHYDE	10.2	
620-23-51-----	M, P-TOLUALDEHYDE	19.0	
66-25-1-----	HEXANAL	9.85	
5779-94-2-----	2,5-Dimethylbenzaldehyde	9.68	
=====	=====	=====	=====

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123008.d

Lab Smp Id: 1838-41-10Client Smp ID: CCV

Inj Date : 23-NOV-2010 10:06

Operator : FMInst ID: hplcf.i

Smp Info : ;1838-41-10;CCV

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 23-Nov-2010 12:36 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1QC Sample: METHSPIKE

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
					ON-COLUMN	FINAL	
	RT	EXP RT	DLT RT	RESPONSE	(ng)	(ug)	
=====	==	=====	=====	=====	=====	=====	
2 FORMALDEHYDE	3.675	3.598	0.077	1127609	10.3002	10.3	
3 ACETALDEHYDE	4.268	4.165	0.103	785758	9.22699	9.23	
4 ACRROLEIN	4.935	4.791	0.144	819165	10.0968	10.1	
5 ACETONE	5.128	4.991	0.137	609926	9.78633	9.79	
6 PROPANAL	5.355	5.198	0.157	579305	9.24166	9.24	
7 CROTONAL	6.221	5.991	0.230	559357	9.82293	9.82	
9 BUTYRALS/MEK	6.835	6.565	0.270	519448	9.59572	9.60	
10 BENZALDEHYDE	7.755	7.405	0.350	367463	9.60466	9.60	
11 ISO-PENTANAL	8.735	8.338	0.397	453422	9.97496	9.97	
12 PENTANAL	9.335	8.898	0.437	433568	9.69947	9.70	
14 O-TOLUALDEHYDE	10.455	9.985	0.470	370488	10.1598	10.2	
15 M,P-TOLUALDEHYDE	10.781	10.338	0.443	582745	19.0230	19.0	
16 HEXANAL	12.415	12.091	0.324	397187	9.84745	9.85	
17 2,5-Dimethylbenzaldehyde	13.041	12.745	0.296	285571	9.68203	9.68	

Air Toxics Ltd.

RECOVERY REPORT

Client Name: Client SDG: 23Nov2010
Sample Matrix: LIQUID Fraction: OTHER
Lab Smp Id: 1838-41-10 Client Smp ID: CCV
Level: MED Operator: FM
Data Type: LC DATA SampleType: METHSPIKE
SpikeList File: 10ug.spk Quant Type: ESTD
Sublist File: all.sub
Method File: /chem/hplcf.i/23Nov2010.b/F10D0914.m
Misc Info:

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	10.0	10.3	103.00	90-110
3 ACETALDEHYDE	10.0	9.23	92.27	90-110
4 ACROLEIN	10.0	10.1	100.97	90-110
5 ACETONE	10.0	9.79	97.86	90-110
6 PROPANAL	10.0	9.24	92.42	90-110
7 CROTONAL	10.0	9.82	98.23	90-110
9 BUTYRALS/MEK	10.0	9.60	95.96	90-110
10 BENZALDEHYDE	10.0	9.60	96.05	90-110
11 ISO-PENTANAL	10.0	9.97	99.75	90-110
12 PENTANAL	10.0	9.70	96.99	90-110
14 O-TOLUALDEHYDE	10.0	10.2	101.60	90-110
15 M,P-TOLUALDEHYDE	20.0	19.0	95.11	90-110
16 HEXANAL	10.0	9.85	98.47	90-110
17 2,5-Dimethylbenzal	10.0	9.68	96.82	90-110

Data File: /chem/hplcf.i/23Nov2010.b/f1123008.d

Page 1

Date : 23-NOV-2010 10:06

Client ID: CCV

Instrument: hplcf.i

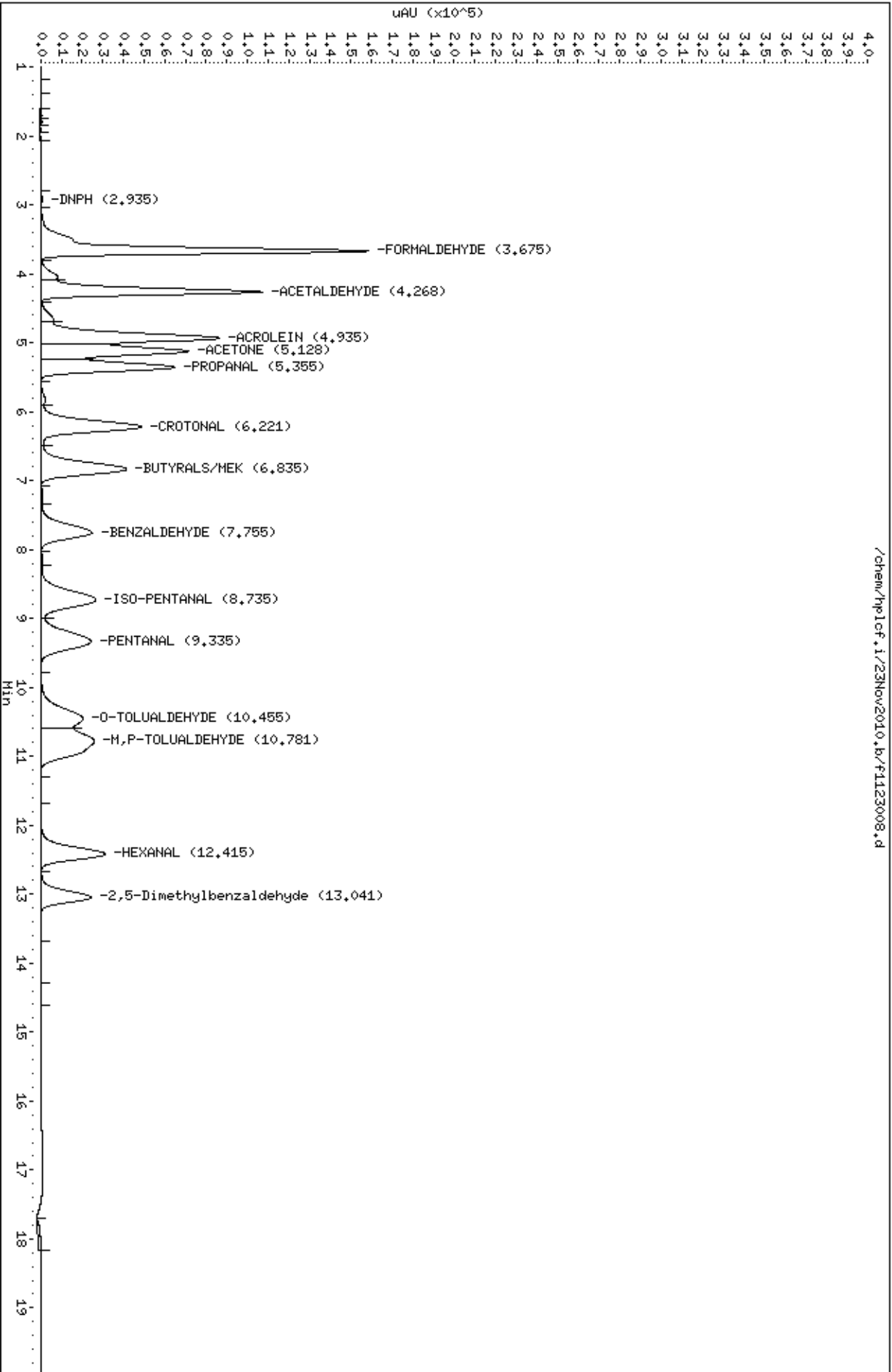
Sample Info: f1838-41-10;CCV

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123009.d

Lab Smp Id: CCBClient Smp ID: LAB BLANK

Inj Date : 23-NOV-2010 10:27

Operator : FMInst ID: hplcf.i

Smp Info : ;CCB;LAB BLANK

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 23-Nov-2010 12:36 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
	RT	EXP RT	DLT RT	RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====		=====	=====	=====
1 DNPH					Compound Not Detected.		
2 FORMALDEHYDE					Compound Not Detected.		
3 ACETALDEHYDE					Compound Not Detected.		
4 ACRYLEIN					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	CONCENTRATIONS					
	RT	EXP	RT	DLT	RT	RESPONSE
=====	==	=====	=====			=====
17 2,5-Dimethylbenzaldehyde						Compound Not Detected.

Data File: /chem/hplcf.i/23Nov2010.b/f1123009.d

Page 1

Date : 23-NOV-2010 10:27

Client ID: LAB BLANK

Instrument: hplcf.i

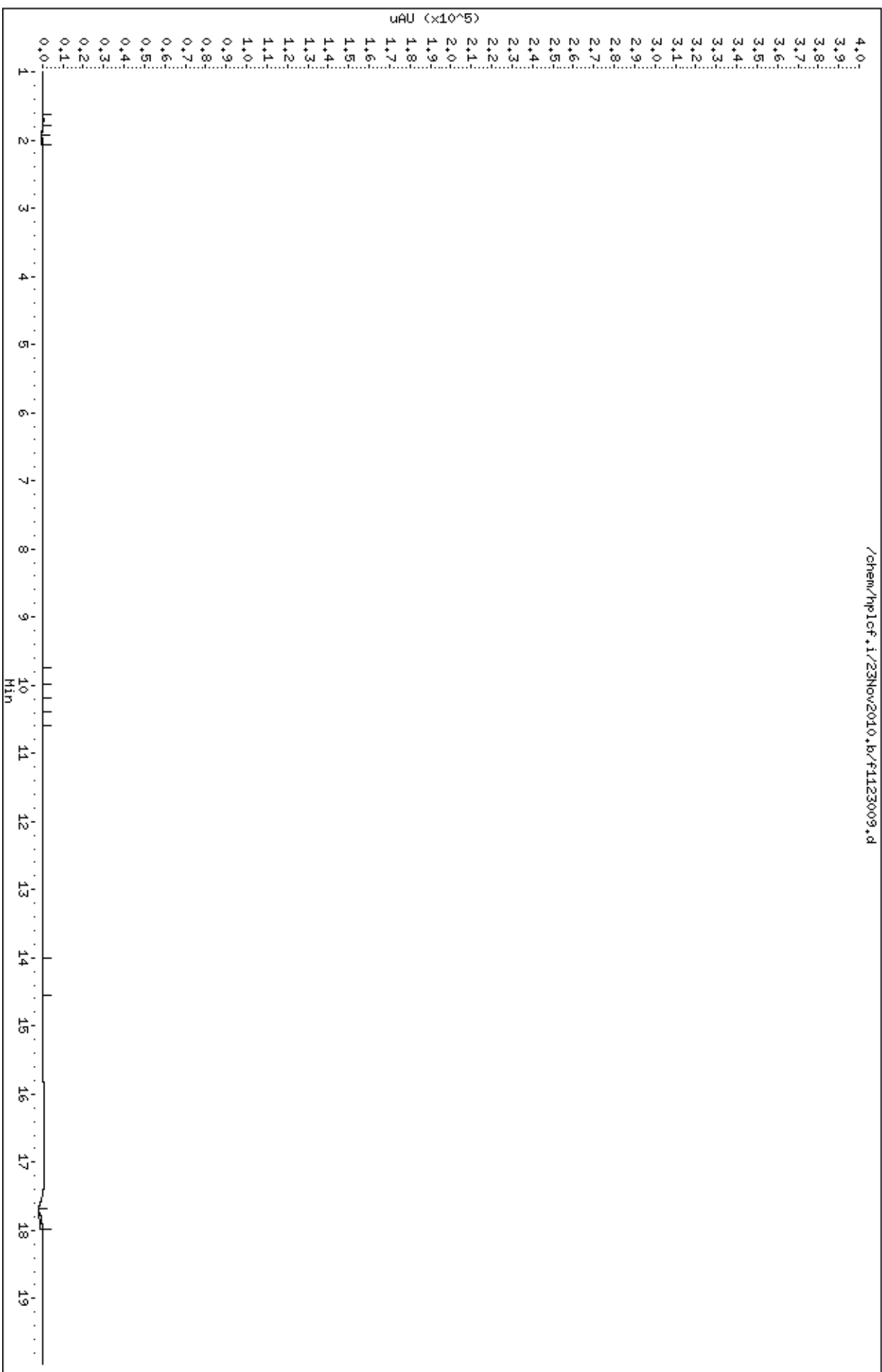
Sample Info: JCCB;LAB BLANK

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

TARGET COMPOUNDS

Client Name:Client SDG: 23Nov2010

Lab Smp Id: 1838-41-10Client Smp ID: CCV

Sample Location:Sample Point:

Sample Date:Date Received:

Sample Matrix: AIRQuant Type: ESTD

Analysis Type: OTHERLevel: MED

Data Type: LC DATAOperator: FM

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	10.1	
107-02-8-----	ACROLEIN	11.0	
67-64-1-----	ACETONE	9.64	
123-38-6-----	PROPANAL	9.13	
123-73-9-----	CROTONAL	10.2	
9999-9999-007--	BUTYRALS/MEK	9.67	
100-52-7-----	BENZALDEHYDE	9.70	
590-86-3-----	ISO-PENTANAL	10.1	
110-62-3-----	PENTANAL	9.66	
529-20-4-----	O-TOLUALDEHYDE	10.4	
620-23-51-----	M, P-TOLUALDEHYDE	18.9	
66-25-1-----	HEXANAL	9.89	
5779-94-2-----	2,5-Dimethylbenzaldehyde	9.67	
=====	=====	=====	=====

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123017.d

Lab Smp Id: 1838-41-10Client Smp ID: CCV

Inj Date : 23-NOV-2010 13:14

Operator : FMInst ID: hplcf.i

Smp Info : ;1838-41-10;CCV

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 23-Nov-2010 12:36 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1QC Sample: METHSPIKE

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
	RT	EXP RT	DLT RT	RESPONSE	ON - COLUMN	FINAL	
					(ng)	(ug)	
=====	==	=====	=====	=====	=====	=====	
2 FORMALDEHYDE	3.673	3.598	0.075	1142002	10.4317	10.4	
3 ACETALDEHYDE	4.273	4.165	0.108	857222	10.0662	10.1	
4 ACROLEIN	4.953	4.791	0.162	891863	10.9928	11.0	
5 ACETONE	5.153	4.991	0.162	600647	9.63745	9.64	
6 PROPANAL	5.380	5.198	0.182	572445	9.13222	9.13	
7 CROTONAL	6.260	5.991	0.269	582083	10.2220	10.2	
9 BUTYRALS/MEK	6.886	6.565	0.321	523307	9.66703	9.67	
10 BENZALDEHYDE	7.819	7.405	0.414	371279	9.70441	9.70	
11 ISO - PENTANAL	8.819	8.338	0.481	459622	10.1113	10.1	
12 PENTANAL	9.440	8.898	0.542	431905	9.66225	9.66	
14 O - TOLUALDEHYDE	10.579	9.985	0.594	379555	10.4084	10.4	
15 M, P - TOLUALDEHYDE	10.899	10.338	0.561	578572	18.8867	18.9	
16 HEXANAL	12.506	12.091	0.415	398772	9.88676	9.89	
17 2,5-Dimethylbenzaldehyde	13.133	12.745	0.388	285282	9.67223	9.67	

Air Toxics Ltd.

RECOVERY REPORT

Client Name: Client SDG: 23Nov2010
Sample Matrix: LIQUID Fraction: OTHER
Lab Smp Id: 1838-41-10 Client Smp ID: CCV
Level: MED Operator: FM
Data Type: LC DATA SampleType: METHSPIKE
SpikeList File: 10ug.spk Quant Type: ESTD
Sublist File: all.sub
Method File: /chem/hplcf.i/23Nov2010.b/F10D0914.m
Misc Info:

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	10.0	10.4	104.32	90-110
3 ACETALDEHYDE	10.0	10.1	100.66	90-110
4 ACROLEIN	10.0	11.0	109.93	90-110
5 ACETONE	10.0	9.64	96.37	90-110
6 PROPANAL	10.0	9.13	91.32	90-110
7 CROTONAL	10.0	10.2	102.22	90-110
9 BUTYRALS/MEK	10.0	9.67	96.67	90-110
10 BENZALDEHYDE	10.0	9.70	97.04	90-110
11 ISO-PENTANAL	10.0	10.1	101.11	90-110
12 PENTANAL	10.0	9.66	96.62	90-110
14 O-TOLUALDEHYDE	10.0	10.4	104.08	90-110
15 M,P-TOLUALDEHYDE	20.0	18.9	94.43	90-110
16 HEXANAL	10.0	9.89	98.87	90-110
17 2,5-Dimethylbenzal	10.0	9.67	96.72	90-110

Data File: /chem/hp1cf.i/23Nov2010.b/f1123017.d

Page 1

Date : 23-NOV-2010 13:14

Client ID: CCV

Instrument: hp1cf.i

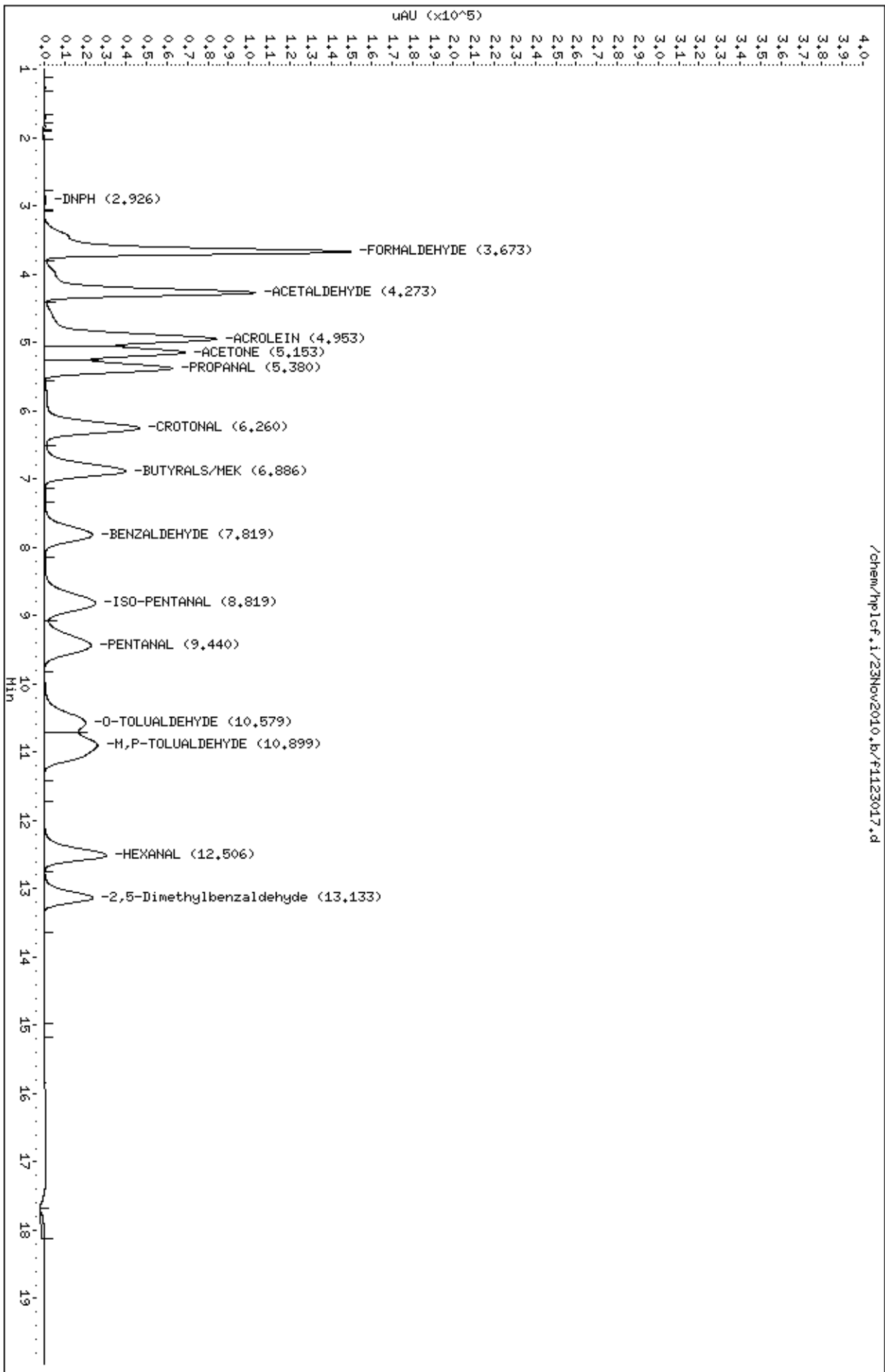
Sample Info: #1838-41-10;CCV

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123018.d

Lab Smp Id: CCBClient Smp ID: LAB BLANK

Inj Date : 23-NOV-2010 13:35

Operator : FMInst ID: hplcf.i

Smp Info : ;CCB;LAB BLANK

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 23-Nov-2010 12:36 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
	RT	EXP RT	DLT RT	RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====		=====	=====	=====
1 DNPH					Compound Not Detected.		
2 FORMALDEHYDE					Compound Not Detected.		
3 ACETALDEHYDE					Compound Not Detected.		
4 ACRROLEIN					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	CONCENTRATIONS					
	RT	EXP RT	DLT	RT	RESPONSE	ON-COLUMN
						FINAL
=====	==	=====	=====		=====	(ng) (ug)
17 2,5-Dimethylbenzaldehyde						=====
					Compound Not Detected.	

Data File: /chem/hp1cf.i/23Nov2010.b/f1123018.d

Page 1

Date : 23-NOV-2010 13:35

Client ID: LAB BLANK

Instrument: hp1cf.i

Sample Info: JCCB;LAB BLANK

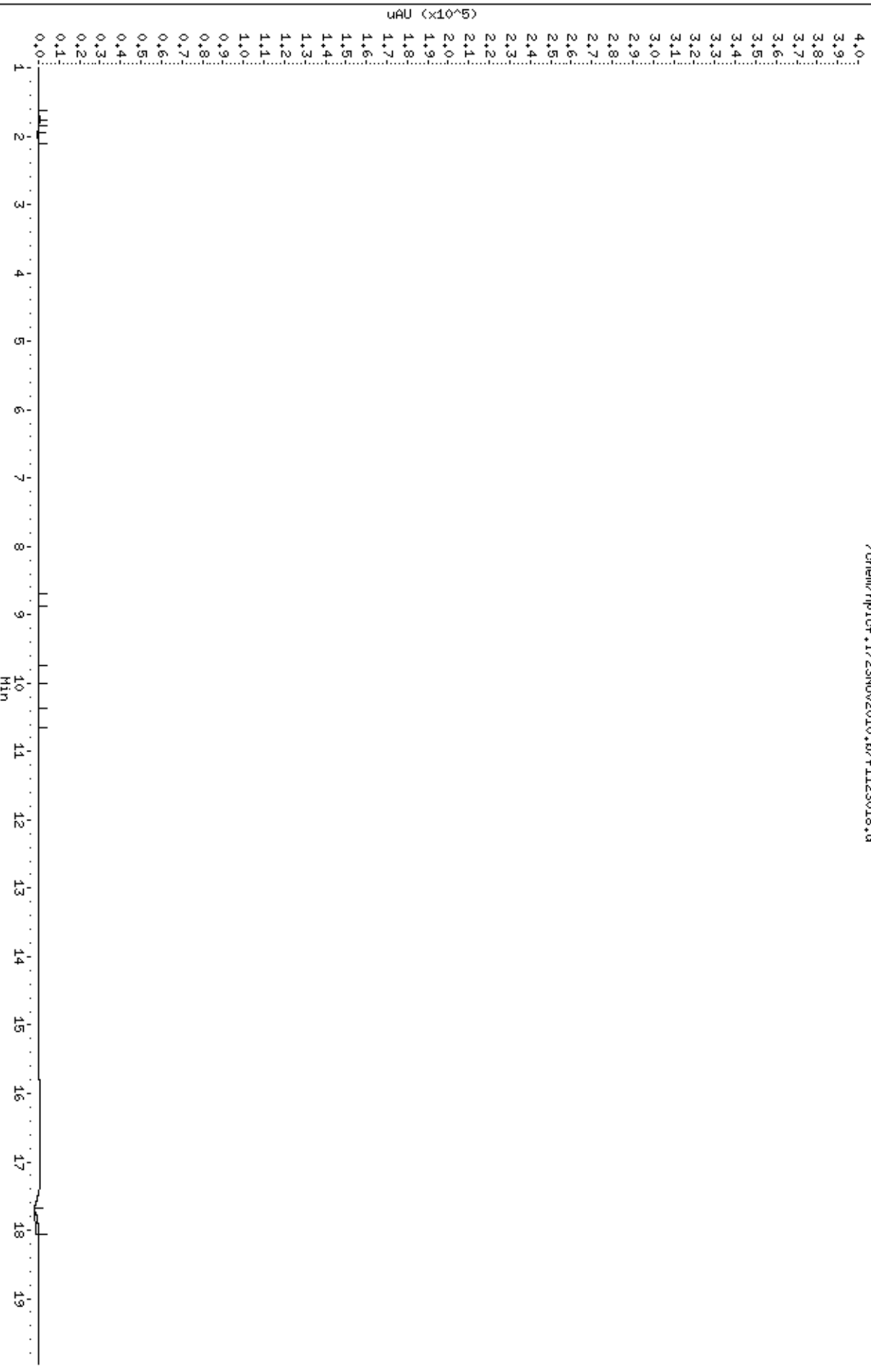
Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00

/chem/hp1cf.i/23Nov2010.b/f1123018.d



Air Toxics Ltd.

TARGET COMPOUNDS

Client Name:Client SDG: 23Nov2010

Lab Smp Id: 1838-41-10Client Smp ID: CCV

Sample Location:Sample Point:

Sample Date:Date Received:

Sample Matrix: AIRQuant Type: ESTD

Analysis Type: OTHERLevel: MED

Data Type: LC DATAOperator: FM

Misc Info:

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/KG) ug	Q
119-26-6-----	DNPH	0.00	
50-00-0-----	FORMALDEHYDE	10.3	
75-07-0-----	ACETALDEHYDE	9.76	
107-02-8-----	ACROLEIN	10.5	
67-64-1-----	ACETONE	9.67	
123-38-6-----	PROPANAL	9.25	
123-73-9-----	CROTONAL	9.91	
9999-9999-007--	BUTYRALS/MEK	9.69	
100-52-7-----	BENZALDEHYDE	9.60	
590-86-3-----	ISO-PENTANAL	9.94	
110-62-3-----	PENTANAL	9.61	
529-20-4-----	O-TOLUALDEHYDE	10.1	
620-23-51-----	M, P-TOLUALDEHYDE	19.2	
66-25-1-----	HEXANAL	9.85	
5779-94-2-----	2,5-Dimethylbenzaldehyde	9.63	
=====		=====	=====

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123030.d

Lab Smp Id: 1838-41-10Client Smp ID: CCV

Inj Date : 23-NOV-2010 17:46

Operator : FMInst ID: hplcf.i

Smp Info : ;1838-41-10;CCV

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 23-Nov-2010 12:36 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1QC Sample: METHSPIKE

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
					ON-COLUMN	FINAL	
	RT	EXP RT	DLT RT	RESPONSE	(ng)	(ug)	
=====	==	=====	=====	=====	=====	=====	
2 FORMALDEHYDE	3.692	3.598	0.094	1128813	10.3112	10.3	
3 ACETALDEHYDE	4.292	4.165	0.127	831255	9.76126	9.76	
4 ACRROLEIN	4.972	4.791	0.181	850078	10.4778	10.5	
5 ACETONE	5.172	4.991	0.181	602540	9.66782	9.67	
6 PROPANAL	5.405	5.198	0.207	579661	9.24734	9.25	
7 CROTONAL	6.285	5.991	0.294	564235	9.90858	9.91	
9 BUTYRALS/MEK	6.905	6.565	0.340	524811	9.69480	9.69	
10 BENZALDEHYDE	7.852	7.405	0.447	367361	9.60199	9.60	
11 ISO-PENTANAL	8.845	8.338	0.507	451853	9.94043	9.94	
12 PENTANAL	9.465	8.898	0.567	429598	9.61065	9.61	
14 O-TOLUALDEHYDE	10.598	9.985	0.613	367745	10.0846	10.1	
15 M,P-TOLUALDEHYDE	10.918	10.338	0.580	588738	19.2186	19.2	
16 HEXANAL	12.525	12.091	0.434	397168	9.84698	9.85	
17 2,5-Dimethylbenzaldehyde	13.145	12.745	0.400	284071	9.63121	9.63	

Air Toxics Ltd.

RECOVERY REPORT

Client Name: Client SDG: 23Nov2010
Sample Matrix: LIQUID Fraction: OTHER
Lab Smp Id: 1838-41-10 Client Smp ID: CCV
Level: MED Operator: FM
Data Type: LC DATA SampleType: METHSPIKE
SpikeList File: 10ug.spk Quant Type: ESTD
Sublist File: all.sub
Method File: /chem/hplcf.i/23Nov2010.b/F10D0914.m
Misc Info:

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	10.0	10.3	103.11	90-110
3 ACETALDEHYDE	10.0	9.76	97.61	90-110
4 ACROLEIN	10.0	10.5	104.78	90-110
5 ACETONE	10.0	9.67	96.68	90-110
6 PROPANAL	10.0	9.25	92.47	90-110
7 CROTONAL	10.0	9.91	99.09	90-110
9 BUTYRALS/MEK	10.0	9.69	96.95	90-110
10 BENZALDEHYDE	10.0	9.60	96.02	90-110
11 ISO-PENTANAL	10.0	9.94	99.40	90-110
12 PENTANAL	10.0	9.61	96.11	90-110
14 O-TOLUALDEHYDE	10.0	10.1	100.85	90-110
15 M,P-TOLUALDEHYDE	20.0	19.2	96.09	90-110
16 HEXANAL	10.0	9.85	98.47	90-110
17 2,5-Dimethylbenzal	10.0	9.63	96.31	90-110

Data File: /chem/hplcf.i/23Nov2010.b/f1123030.d

Page 1

Date : 23-NOV-2010 17:46

Client ID: CCV

Instrument: hplcf.i

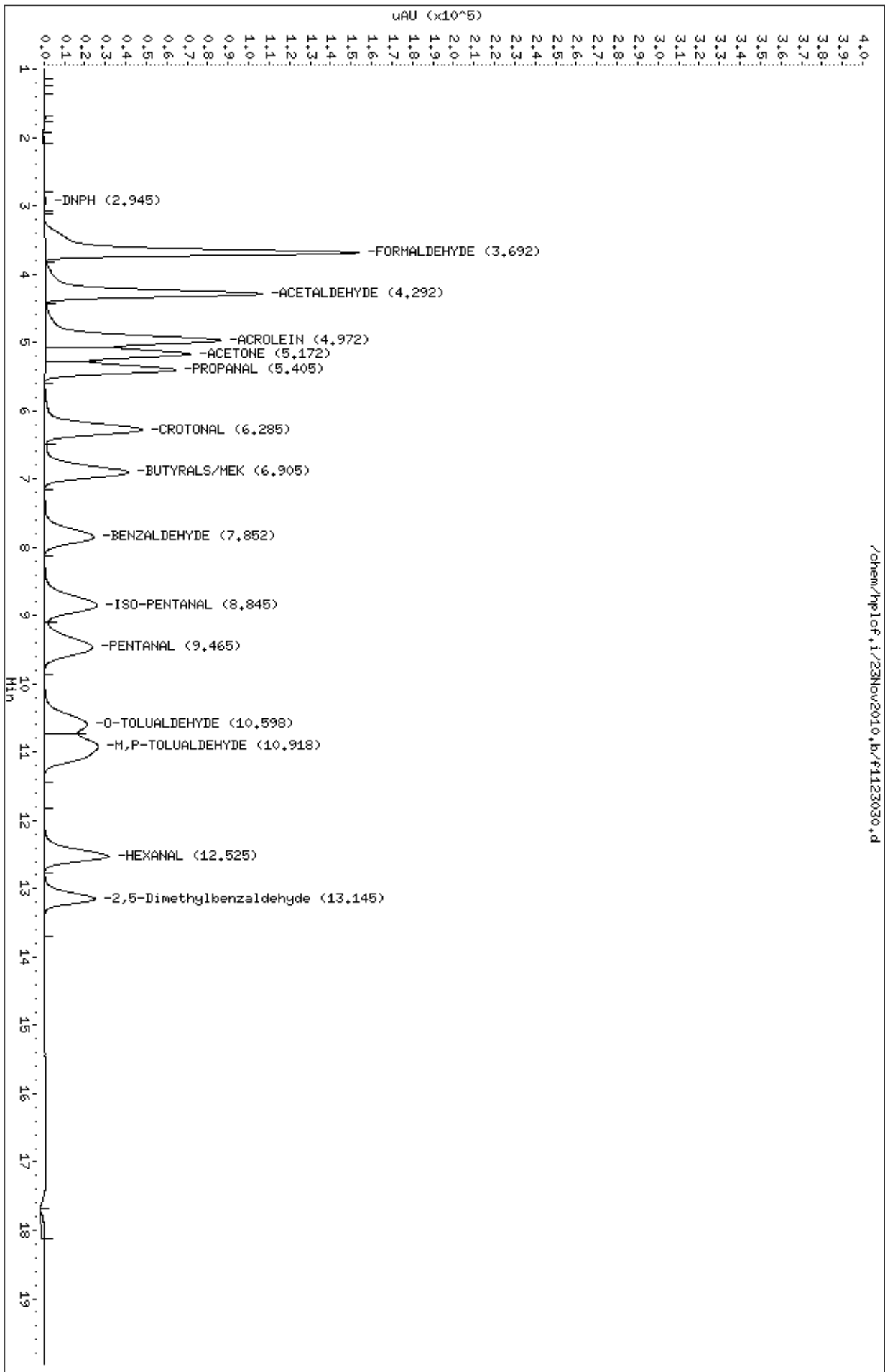
Sample Info: #1838-41-10;CCV

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123031.d

Lab Smp Id: CCBClient Smp ID: LAB BLANK

Inj Date : 23-NOV-2010 18:06

Operator : FMInst ID: hplcf.i

Smp Info : ;CCB;LAB BLANK

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 23-Nov-2010 12:36 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)	
=====	==	=====	=====	=====	=====	=====	
1 DNPH				Compound Not Detected.			
2 FORMALDEHYDE				Compound Not Detected.			
3 ACETALDEHYDE				Compound Not Detected.			
4 ACRROLEIN				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	CONCENTRATIONS					
	RT	EXP RT	DLT	RT	RESPONSE	ON-COLUMN
						FINAL
=====	==	=====	=====		=====	(ng) (ug)
17 2,5-Dimethylbenzaldehyde						=====
					Compound Not Detected.	

Data File: /chem/hplcf.i/23Nov2010.b/f1123031.d

Page 1

Date : 23-NOV-2010 18:06

Client ID: LAB BLANK

Instrument: hplcf.i

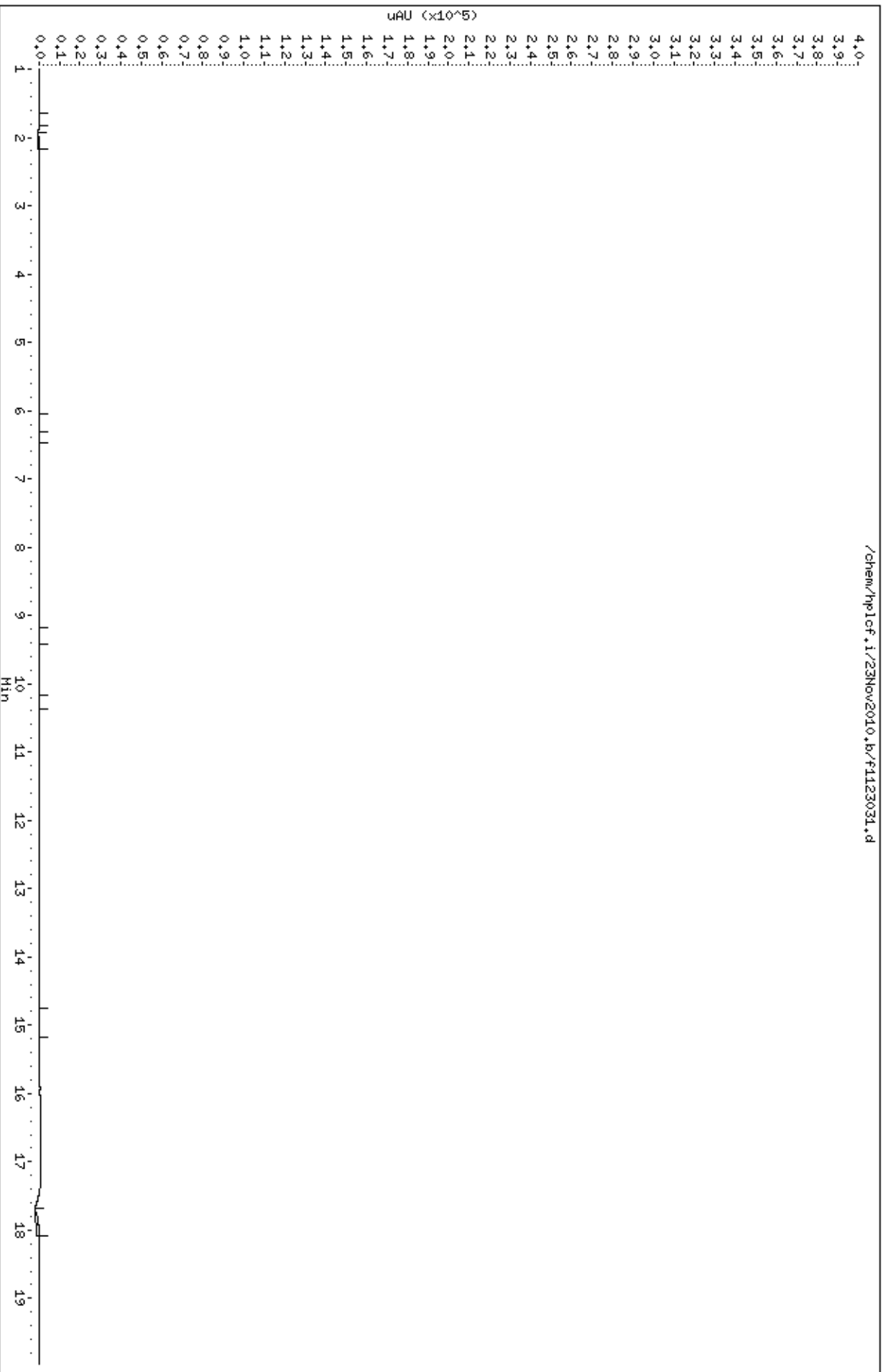
Sample Info: JCCB;LAB BLANK

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



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: 23Nov2010
Client Smp ID: CCV
Sample Point:
Date Received:
Quant Type: ESTD
Level: MED
Operator: FM

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	11.0	
67-64-1-----	ACETONE	9.73	
123-38-6-----	PROPANAL	9.14	
123-73-9-----	CROTONAL	10.3	
9999-9999-007--	BUTYRALS/MEK	9.66	
100-52-7-----	BENZALDEHYDE	9.72	
590-86-3-----	ISO-PENTANAL	10.1	
110-62-3-----	PENTANAL	9.63	
529-20-4-----	O-TOLUALDEHYDE	10.3	
620-23-51-----	M, P-TOLUALDEHYDE	18.8	
66-25-1-----	HEXANAL	10.1	
5779-94-2-----	2,5-Dimethylbenzaldehyde	9.67	
=====	=====	=====	=====

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123042.d

Lab Smp Id: 1838-41-10Client Smp ID: CCV

Inj Date : 23-NOV-2010 21:56

Operator : FMInst ID: hplcf.i

Smp Info : ;1838-41-10;CCV

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1QC Sample: METHSPIKE

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS					
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN	FINAL
					(ng)	(ug)
=====	==	=====	=====	=====	=====	=====
2 FORMALDEHYDE	3.651	3.598	0.053	1145512	10.4638	10.5
3 ACETALDEHYDE	4.245	4.165	0.080	860273	10.1020	10.1
4 ACROLEIN	4.911	4.791	0.120	891474	10.9880	11.0
5 ACETONE	5.105	4.991	0.114	606562	9.73236	9.73
6 PROPANAL	5.331	5.198	0.133	573213	9.14448	9.14
7 CROTONAL	6.185	5.991	0.194	585724	10.2859	10.3
9 BUTYRALS/MEK	6.805	6.565	0.240	522783	9.65734	9.66
10 BENZALDEHYDE	7.711	7.405	0.306	372065	9.72494	9.72
11 ISO-PENTANAL	8.698	8.338	0.360	459538	10.1095	10.1
12 PENTANAL	9.298	8.898	0.400	430288	9.62608	9.63
14 O-TOLUALDEHYDE	10.425	9.985	0.440	376945	10.3369	10.3
15 M,P-TOLUALDEHYDE	10.751	10.338	0.413	577449	18.8501	18.8
16 HEXANAL	12.405	12.091	0.314	406661	10.0823	10.1
17 2,5-Dimethylbenzaldehyde	13.038	12.745	0.293	285337	9.67411	9.67

Air Toxics Ltd.

RECOVERY REPORT

Client Name: Client SDG: 23Nov2010
Sample Matrix: LIQUID Fraction: OTHER
Lab Smp Id: 1838-41-10 Client Smp ID: CCV
Level: MED Operator: FM
Data Type: LC DATA SampleType: METHSPIKE
SpikeList File: 10ug.spk Quant Type: ESTD
Sublist File: all.sub
Method File: /chem/hplcf.i/23Nov2010.b/F10D0914.m
Misc Info:

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	10.0	10.5	104.64	90-110
3 ACETALDEHYDE	10.0	10.1	101.02	90-110
4 ACROLEIN	10.0	11.0	109.88	90-110
5 ACETONE	10.0	9.73	97.32	90-110
6 PROPANAL	10.0	9.14	91.44	90-110
7 CROTONAL	10.0	10.3	102.86	90-110
9 BUTYRALS/MEK	10.0	9.66	96.57	90-110
10 BENZALDEHYDE	10.0	9.72	97.25	90-110
11 ISO-PENTANAL	10.0	10.1	101.09	90-110
12 PENTANAL	10.0	9.63	96.26	90-110
14 O-TOLUALDEHYDE	10.0	10.3	103.37	90-110
15 M,P-TOLUALDEHYDE	20.0	18.8	94.25	90-110
16 HEXANAL	10.0	10.1	100.82	90-110
17 2,5-Dimethylbenzal	10.0	9.67	96.74	90-110

Data File: /chem/hplcf,i/23Nov2010,b/f1123042.d

Date : 23-NOV-2010 21:56

Client ID: CCV

Sample Info: f1838-41-10;CCV

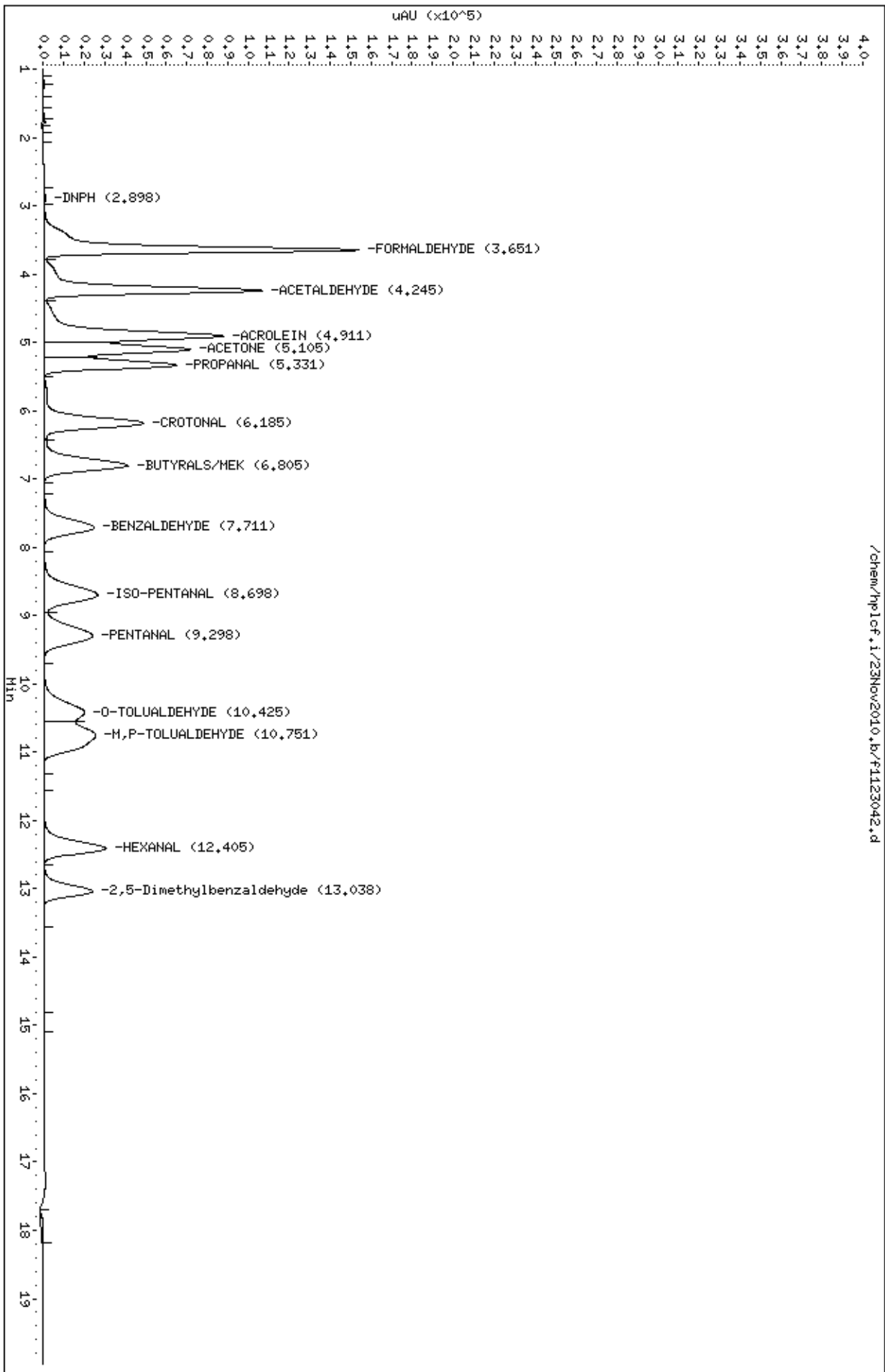
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hplcf,i

Operator: FM

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123043.d

Lab Smp Id: CCBClient Smp ID: LAB BLANK

Inj Date : 23-NOV-2010 22:17

Operator : FMInst ID: hplcf.i

Smp Info : ;CCB;LAB BLANK

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
	RT	EXP RT	DLT RT	RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====		=====	=====	=====
1 DNPH					Compound Not Detected.		
2 FORMALDEHYDE					Compound Not Detected.		
3 ACETALDEHYDE					Compound Not Detected.		
4 ACRROLEIN					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	CONCENTRATIONS					
	RT	EXP RT	DLT	RT	RESPONSE	ON-COLUMN
						FINAL
=====	==	=====	=====		=====	(ng) (ug)
17 2,5-Dimethylbenzaldehyde						=====
					Compound Not Detected.	

Data File: /chem/hplcf.i/23Nov2010.b/f1123043.d

Page 1

Date : 23-NOV-2010 22:17

Client ID: LAB BLANK

Instrument: hplcf.i

Sample Info: JCCB;LAB BLANK

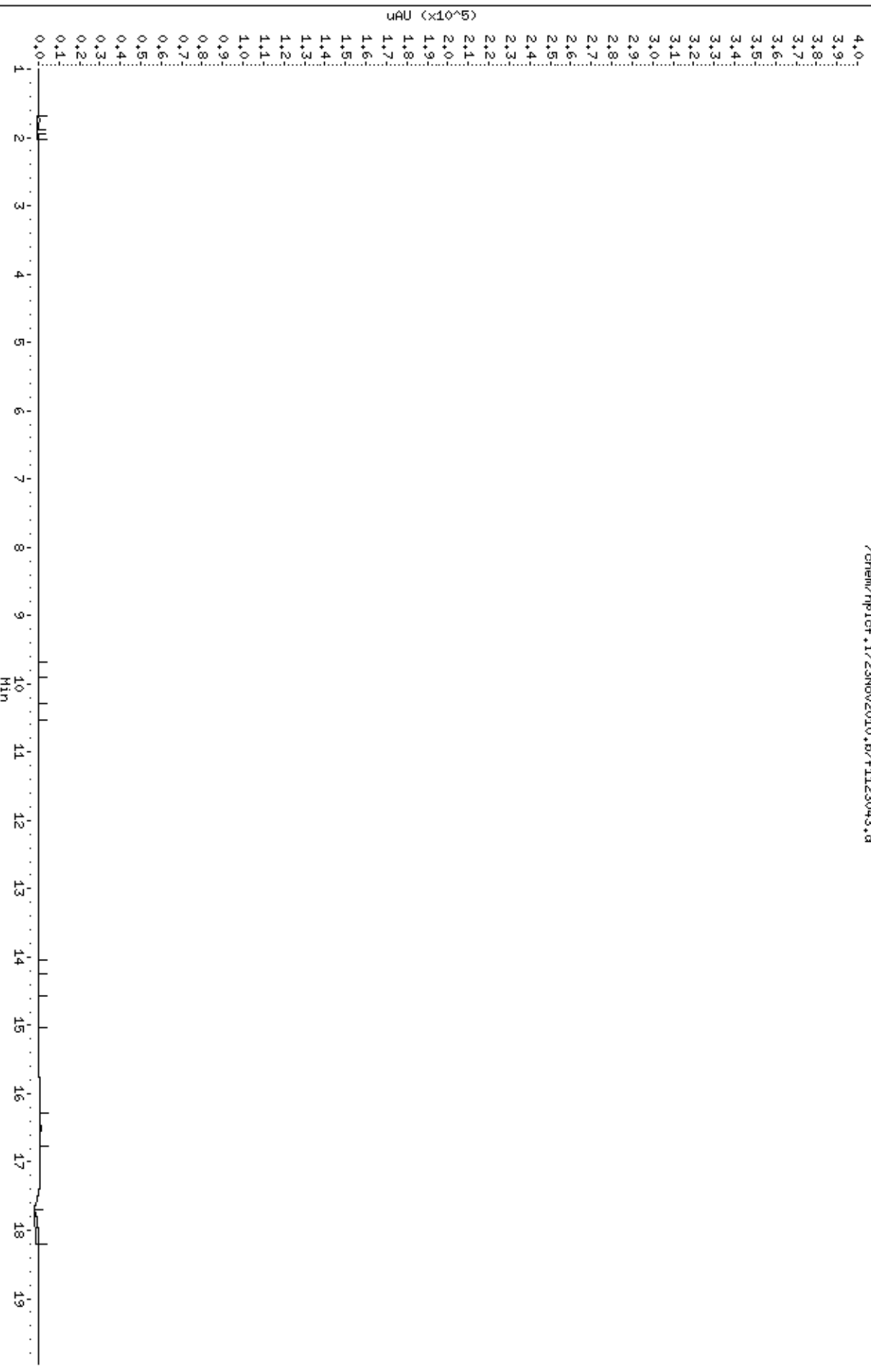
Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00

/chem/hplcf.i/23Nov2010.b/f1123043.d



Air Toxics Ltd.

TARGET COMPOUNDS

Client Name:Client SDG: 23Nov2010

Lab Smp Id: 1838-41-10Client Smp ID: CCV

Sample Location:Sample Point:

Sample Date:Date Received:

Sample Matrix: AIRQuant Type: ESTD

Analysis Type: OTHERLevel: MED

Data Type: LC DATAOperator: FM

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	10.1	
107-02-8-----	ACROLEIN	11.0	
67-64-1-----	ACETONE	9.70	
123-38-6-----	PROPANAL	9.04	
123-73-9-----	CROTONAL	9.95	
9999-9999-007--	BUTYRALS/MEK	9.60	
100-52-7-----	BENZALDEHYDE	9.77	
590-86-3-----	ISO-PENTANAL	10.0	
110-62-3-----	PENTANAL	9.58	
529-20-4-----	O-TOLUALDEHYDE	10.4	
620-23-51-----	M, P-TOLUALDEHYDE	18.6	
66-25-1-----	HEXANAL	10.0	
5779-94-2-----	2,5-Dimethylbenzaldehyde	9.77	
=====	=====	=====	=====

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123054.d

Lab Smp Id: 1838-41-10Client Smp ID: CCV

Inj Date : 24-NOV-2010 02:07

Operator : FMInst ID: hplcf.i

Smp Info : ;1838-41-10;CCV

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1QC Sample: METHSPIKE

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
					ON-COLUMN	FINAL	
	RT	EXP RT	DLT RT	RESPONSE	(ng)	(ug)	
=====	==	=====	=====	=====	=====	=====	
2 FORMALDEHYDE	3.664	3.598	0.066	1143629	10.4466	10.4	
3 ACETALDEHYDE	4.257	4.165	0.092	857356	10.0678	10.1	
4 ACROLEIN	4.917	4.791	0.126	894325	11.0232	11.0(R)	
5 ACETONE	5.111	4.991	0.120	604454	9.69854	9.70	
6 PROPANAL	5.337	5.198	0.139	567015	9.04560	9.04	
7 CROTONAL	6.184	5.991	0.193	566512	9.94857	9.95	
9 BUTYRALS/MEK	6.790	6.565	0.225	519531	9.59726	9.60	
10 BENZALDEHYDE	7.691	7.405	0.286	373685	9.76728	9.77	
11 ISO-PENTANAL	8.657	8.338	0.319	457136	10.0567	10.0	
12 PENTANAL	9.251	8.898	0.353	428152	9.57831	9.58	
14 O-TOLUALDEHYDE	10.364	9.985	0.379	379507	10.4071	10.4	
15 M,P-TOLUALDEHYDE	10.691	10.338	0.353	570305	18.6169	18.6	
16 HEXANAL	12.357	12.091	0.266	404881	10.0382	10.0	
17 2,5-Dimethylbenzaldehyde	12.984	12.745	0.239	288205	9.77135	9.77	

QC Flag Legend

R - Spike/Surrogate failed recovery limits.

Air Toxics Ltd.

RECOVERY REPORT

Client Name: Client SDG: 23Nov2010
Sample Matrix: LIQUID Fraction: OTHER
Lab Smp Id: 1838-41-10 Client Smp ID: CCV
Level: MED Operator: FM
Data Type: LC DATA SampleType: METHSPIKE
SpikeList File: 10ug.spk Quant Type: ESTD
Sublist File: all.sub
Method File: /chem/hplcf.i/23Nov2010.b/F10D0914.m
Misc Info:

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	10.0	10.4	104.47	90-110
3 ACETALDEHYDE	10.0	10.1	100.68	90-110
4 ACROLEIN	10.0	11.0	110.23*	90-110
5 ACETONE	10.0	9.70	96.99	90-110
6 PROPANAL	10.0	9.04	90.46	90-110
7 CROTONAL	10.0	9.95	99.49	90-110
9 BUTYRALS/MEK	10.0	9.60	95.97	90-110
10 BENZALDEHYDE	10.0	9.77	97.67	90-110
11 ISO-PENTANAL	10.0	10.0	100.57	90-110
12 PENTANAL	10.0	9.58	95.78	90-110
14 O-TOLUALDEHYDE	10.0	10.4	104.07	90-110
15 M,P-TOLUALDEHYDE	20.0	18.6	93.08	90-110
16 HEXANAL	10.0	10.0	100.38	90-110
17 2,5-Dimethylbenzal	10.0	9.77	97.71	90-110

Data File: /chem/hp1cf.i/23Nov2010.b/f1123054.d

Page 1

Date : 24-NOV-2010 02:07

Client ID: CCV

Instrument: hp1cf.i

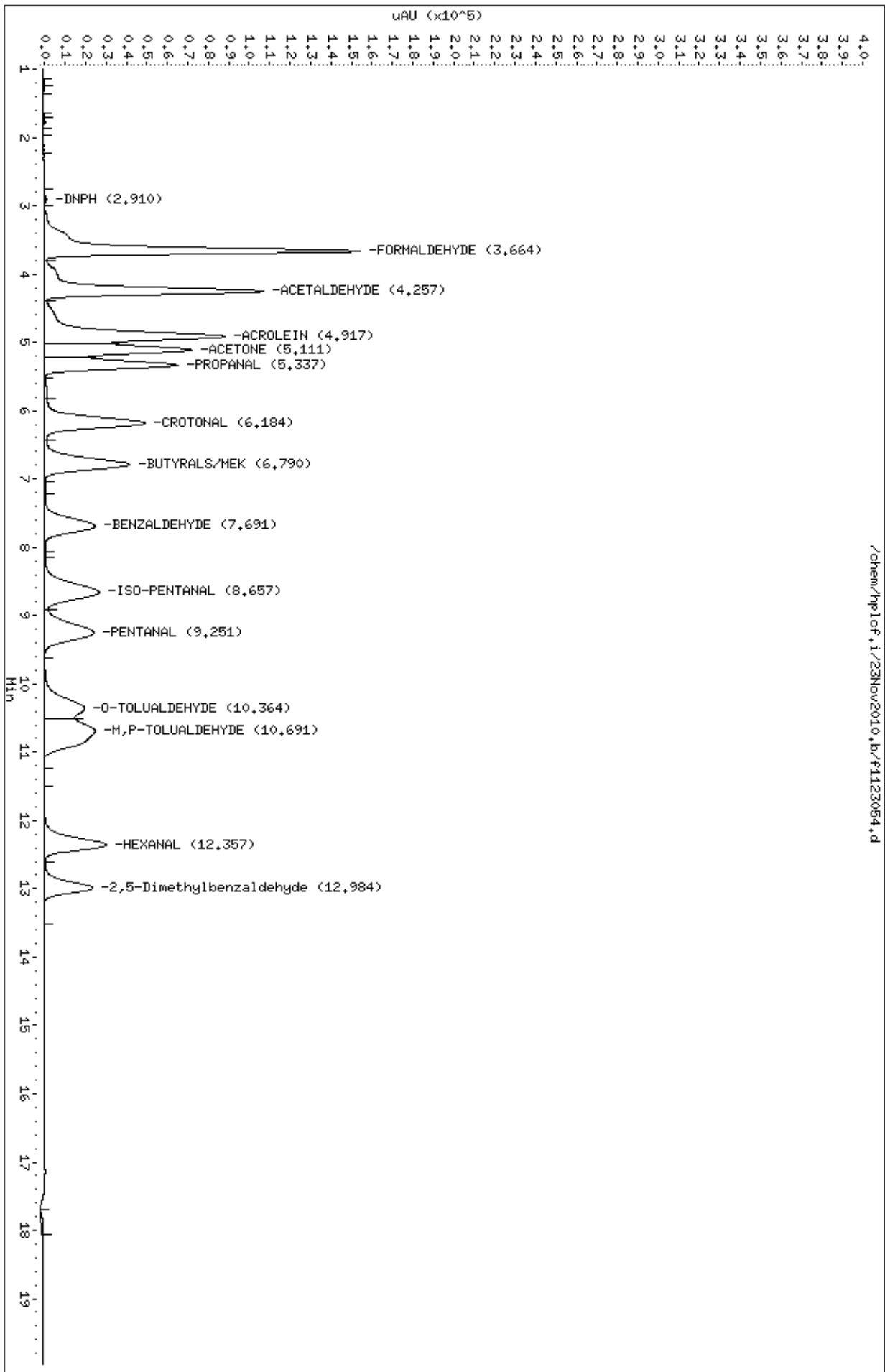
Sample Info: #1838-41-10;CCV

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123055.d

Lab Smp Id: CCBClient Smp ID: LAB BLANK

Inj Date : 24-NOV-2010 02:28

Operator : FMInst ID: hplcf.i

Smp Info : ;CCB;LAB BLANK

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
	RT	EXP RT	DLT RT	RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)
=====	==	=====	=====		=====	=====	=====
1 DNPH					Compound Not Detected.		
2 FORMALDEHYDE					Compound Not Detected.		
3 ACETALDEHYDE					Compound Not Detected.		
4 ACRROLEIN					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	FINAL
							(ng)	(ug)
=====	==	=====	=====			=====	=====	
17 2,5-Dimethylbenzaldehyde						Compound Not Detected.		

Data File: /chem/hplcf.i/23Nov2010.b/f1123055.d

Page 1

Date : 24-NOV-2010 02:28

Client ID: LAB BLANK

Instrument: hplcf.i

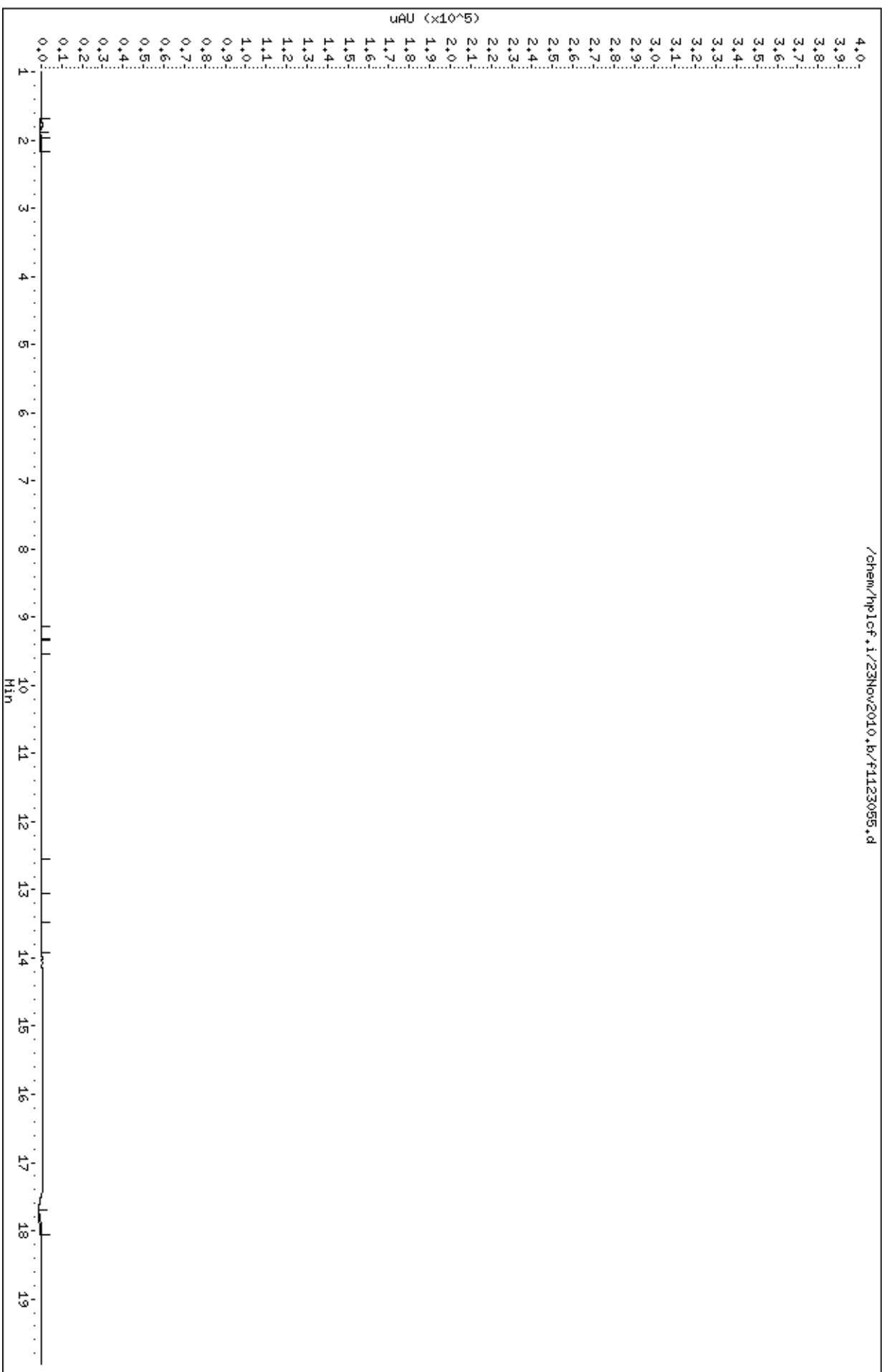
Sample Info: JCCB;LAB BLANK

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



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: 23Nov2010

Client Smp ID: CCV

Sample Point:

Date Received:

Quant Type: ESTD

Level: MED

Operator: FM

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/KG) ug	Q
119-26-6-----	DNPH	0.00	
50-00-0-----	FORMALDEHYDE	10.2	
75-07-0-----	ACETALDEHYDE	10.0	
107-02-8-----	ACROLEIN	10.7	
67-64-1-----	ACETONE	9.71	
123-38-6-----	PROPANAL	9.38	
123-73-9-----	CROTONAL	9.90	
9999-9999-007--	BUTYRALS/MEK	9.83	
100-52-7-----	BENZALDEHYDE	9.80	
590-86-3-----	ISO-PENTANAL	10.1	
110-62-3-----	PENTANAL	9.76	
529-20-4-----	O-TOLUALDEHYDE	10.1	
620-23-51-----	M, P-TOLUALDEHYDE	19.2	
66-25-1-----	HEXANAL	9.98	
5779-94-2-----	2,5-Dimethylbenzaldehyde	9.68	
=====	=====	=====	=====

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123066.d

Lab Smp Id: 1838-41-10Client Smp ID: CCV

Inj Date : 24-NOV-2010 06:17

Operator : FMInst ID: hplcf.i

Smp Info : ;1838-41-10;CCV

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 24-Nov-2010 07:34 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1QC Sample: METHSPIKE

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
					ON-COLUMN	FINAL	
	RT	EXP RT	DLT RT	RESPONSE	(ng)	(ug)	
=====	==	=====	=====	=====	=====	=====	
2 FORMALDEHYDE	3.603	3.598	0.005	1123159	10.2596	10.2	
3 ACETALDEHYDE	4.183	4.165	0.018	856046	10.0524	10.0	
4 ACROLEIN	4.836	4.791	0.045	868344	10.7029	10.7	
5 ACETONE	5.030	4.991	0.039	605339	9.71274	9.71	
6 PROPANAL	5.250	5.198	0.052	588193	9.38346	9.38	
7 CROTONAL	6.090	5.991	0.099	563811	9.90114	9.90	
9 BUTYRALS/MEK	6.690	6.565	0.125	532279	9.83275	9.83	
10 BENZALDEHYDE	7.577	7.405	0.172	375085	9.80388	9.80	
11 ISO-PENTANAL	8.537	8.338	0.199	457419	10.0629	10.1	
12 PENTANAL	9.123	8.898	0.225	436465	9.76427	9.76	
14 O-TOLUALDEHYDE	10.236	9.985	0.251	369668	10.1373	10.1	
15 M,P-TOLUALDEHYDE	10.583	10.338	0.245	588967	19.2261	19.2	
16 HEXANAL	12.277	12.091	0.186	402594	9.98151	9.98	
17 2,5-Dimethylbenzaldehyde	12.916	12.745	0.171	285453	9.67805	9.68	

Air Toxics Ltd.

RECOVERY REPORT

Client Name: Client SDG: 23Nov2010
Sample Matrix: LIQUID Fraction: OTHER
Lab Smp Id: 1838-41-10 Client Smp ID: CCV
Level: MED Operator: FM
Data Type: LC DATA SampleType: METHSPIKE
SpikeList File: 10ug.spk Quant Type: ESTD
Sublist File: all.sub
Method File: /chem/hplcf.i/23Nov2010.b/F10D0914.m
Misc Info:

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	10.0	10.2	102.60	90-110
3 ACETALDEHYDE	10.0	10.0	100.52	90-110
4 ACROLEIN	10.0	10.7	107.03	90-110
5 ACETONE	10.0	9.71	97.13	90-110
6 PROPANAL	10.0	9.38	93.83	90-110
7 CROTONAL	10.0	9.90	99.01	90-110
9 BUTYRALS/MEK	10.0	9.83	98.33	90-110
10 BENZALDEHYDE	10.0	9.80	98.04	90-110
11 ISO-PENTANAL	10.0	10.1	100.63	90-110
12 PENTANAL	10.0	9.76	97.64	90-110
14 O-TOLUALDEHYDE	10.0	10.1	101.37	90-110
15 M,P-TOLUALDEHYDE	20.0	19.2	96.13	90-110
16 HEXANAL	10.0	9.98	99.82	90-110
17 2,5-Dimethylbenzal	10.0	9.68	96.78	90-110

Data File: /chem/hplcf.i/23Nov2010.b/f1123066.d

Date : 24-NOV-2010 06:17

Client ID: CCV

Sample Info: f1838-41-10;CCV

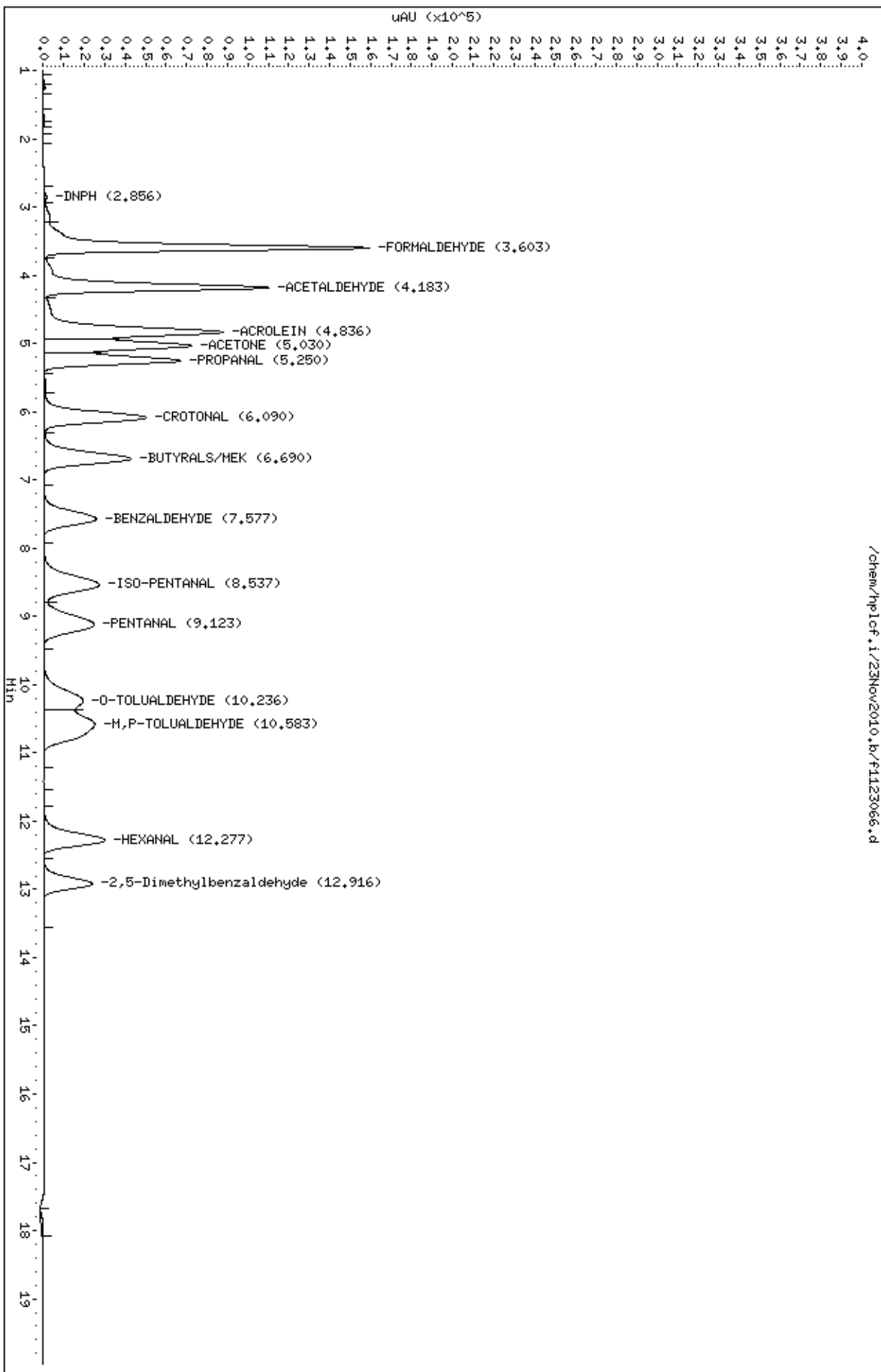
Purge Volume: 5.0

Column phase: C-18 REVERSE PHASE

Instrument: hplcf.i

Operator: FM

Column diameter: 2.00



Client Sample ID: LCS

Lab ID#: 1011418A-17A

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123019
Dil. Factor: 1.00

Date of Collection: NA
Date of Analysis: 11/23/10 01:56 PM
Date of Extraction: 11/23/10

Compound	%Recovery
Formaldehyde	99

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

RECOVERY REPORT

Client Name:	Client SDG: 23Nov2010
Sample Matrix: LIQUID	Fraction: OTHER
Lab Smp Id: 1011418A	Client Smp ID: LCS
Level: MED	Operator: FM
Data Type: LC DATA	SampleType: LCS
SpikeList File: Form15T011.spk	Quant Type: ESTD
Sublist File: form.sub	
Method File: /chem/hplcf.i/23Nov2010.b/F10D0914.m	
Misc Info:	

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	15.0	14.9	99.21	85-115

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123019.d

Lab Smp Id: 1011418AClient Smp ID: LCS

Inj Date : 23-NOV-2010 13:56

Operator : FMInst ID: hplcf.i

Smp Info : ;1011418A;LCS

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 23-Nov-2010 12:36 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1QC Sample: LCS

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)	
=====	==	=====	=====	=====	=====	=====	
1 DNPH	2.905	2.885	0.020	26210508			
2 FORMALDEHYDE	3.684	3.598	0.086	4072665	37.2020	14.9	

Data File: /chem/hp1cf.i/23Nov2010.b/f1123019.d

Page 1

Date : 23-NOV-2010 13:56

Client ID: LCS

Instrument: hp1cf.i

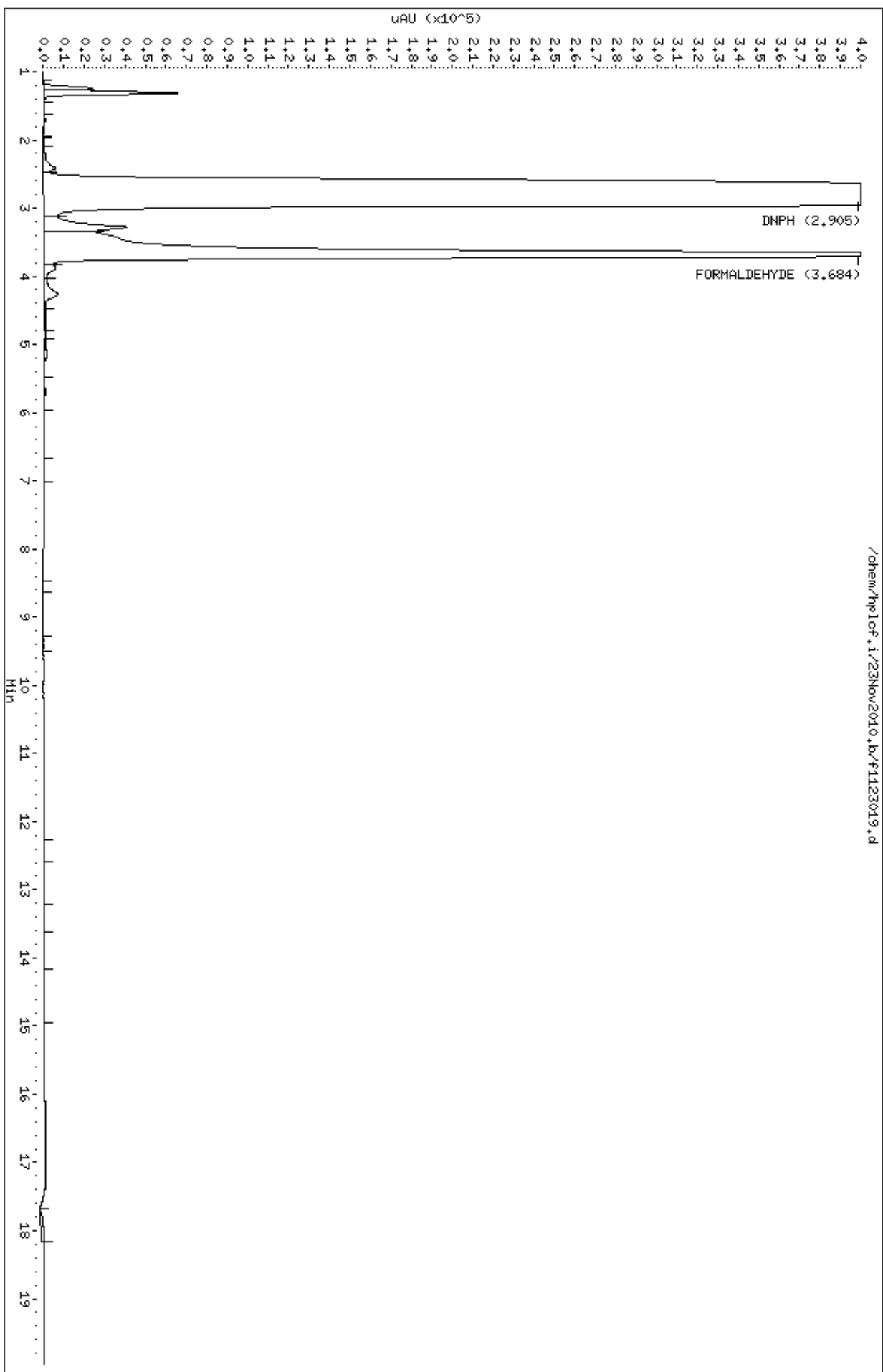
Sample Info: #101141801;LCS

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



Client Sample ID: LCSD

Lab ID#: 1011418A-17AA

RADIELLO ANALYSIS BY HPLC/UV

File Name: f1123020
Dil. Factor: 1.00

Date of Collection: NA
Date of Analysis: 11/23/10 02:17 PM
Date of Extraction: 11/23/10

Compound	%Recovery
Formaldehyde	99

Container Type: Radiello 165 (Aldehydes)

Air Toxics Ltd.

RECOVERY REPORT

Client Name:
Sample Matrix: LIQUID
Lab Smp Id: 1011418A
Level: MED
Data Type: LC DATA
SpikeList File: Form15T011.spk
Sublist File: form.sub
Method File: /chem/hplcf.i/23Nov2010.b/F10D0914.m
Misc Info:

Client SDG: 23Nov2010
Fraction: OTHER
Client Smp ID: LCSD
Operator: FM
SampleType: LCSD
Quant Type: ESTD

SPIKE COMPOUND	CONC ADDED ug	CONC RECOVERED ug	% RECOVERED	LIMITS
2 FORMALDEHYDE	15.0	14.8	98.91	85-115

Air Toxics Ltd.

Aldehydes/Ketones To-11,TO-5,CARB430,0011, Radiello 165

Data file : /chem/hplcf.i/23Nov2010.b/f1123020.d

Lab Smp Id: 1011418AClient Smp ID: LCSD

Inj Date : 23-NOV-2010 14:17

Operator : FMInst ID: hplcf.i

Smp Info : ;1011418A;LCSD

Misc Info :

Comment :

Method : /chem/hplcf.i/23Nov2010.b/F10D0914.m

Meth Date : 23-Nov-2010 12:36 cleafQuant Type: ESTD

Cal Date : 14-SEP-2010 18:30Cal File: f0914014.d

Als bottle: 1QC Sample: LCSD

Dil Factor: 1.00000

Integrator: FalconCompound 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 VariableLocal Compound Variable

Compounds	CONCENTRATIONS						
	RT	EXP RT	DLT RT	RESPONSE	ON-COLUMN (ng)	FINAL (ug)	
=====	==	=====	=====	=====	=====	=====	
1 DNPH	2.898	2.885	0.013	23496381			
2 FORMALDEHYDE	3.665	3.598	0.067	4060628	37.0921	14.8	

Data File: /chem/hp1cf.i/23Nov2010.b/f1123020.d

Page 1

Date : 23-NOV-2010 14:17

Client ID: LCSD

Instrument: hp1cf.i

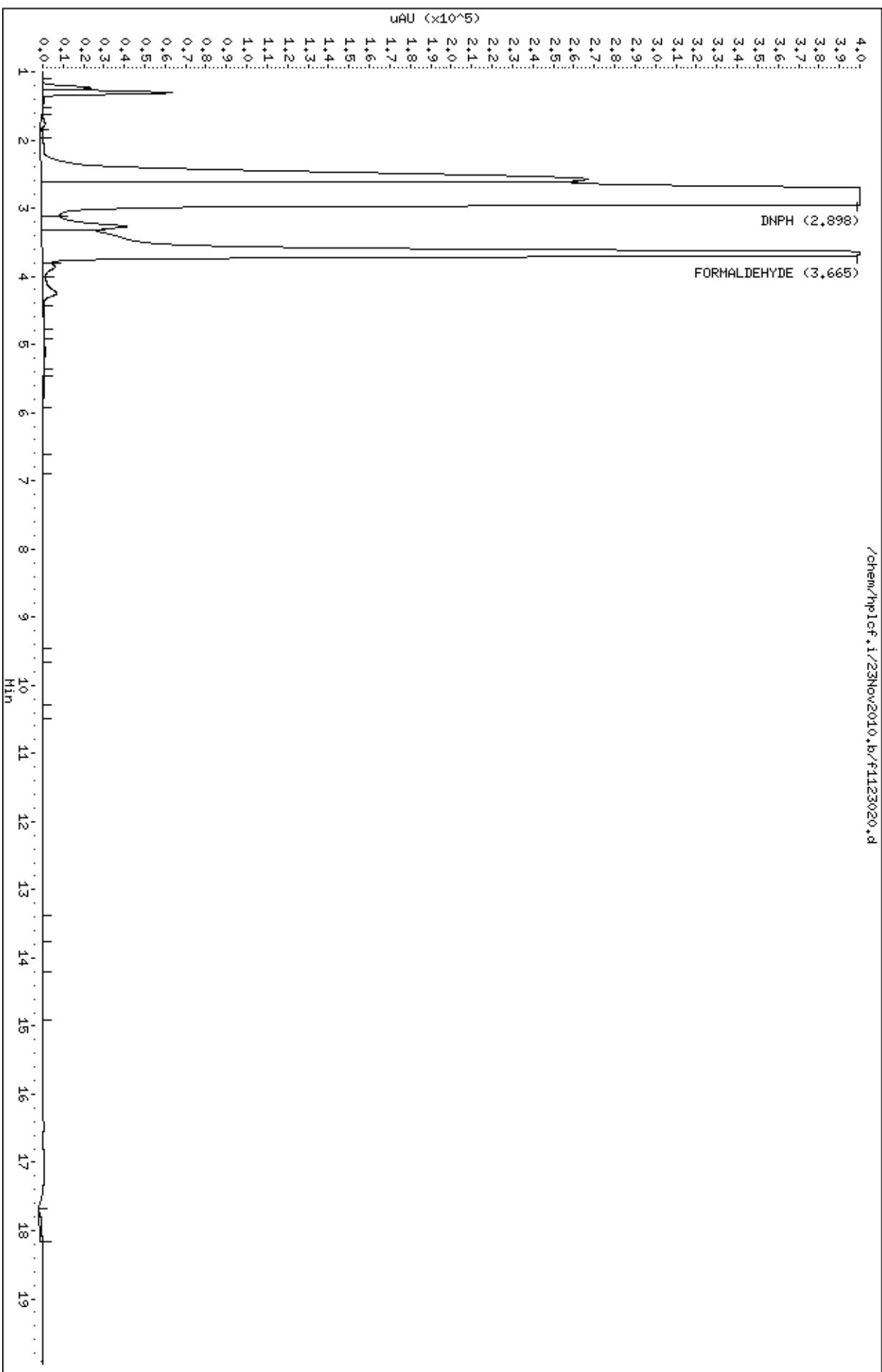
Sample Info: #10114180;LCSD

Purge Volume: 5.0

Operator: FM

Column phase: C-18 REVERSE PHASE

Column diameter: 2.00



TO-11 Extraction Logbook

@Air Toxics Ltd.

Log Book #: 2010

ATL Work Order # 1011418A

Final Solvent: ACN

Solvent Lot #: D8257

Date Extracted 11/23/10

Spike I.D.: 1476-1963-500ug

Spiked By: Fm

Spike Witness: CM

ATL Fraction	Clients I.D. #	Cartridge type & size	Final Volume (ml)	Spike Amount (ul)	Comments
1011418A BLK	118683	Rad 165 10204	2.0	NA	
2CS	4	10048		30	
-01A	5 118683	NA		NA	
-02A	6 4				
-03A	7 5				
-04A	118699 6				
-05A	118700 7				
-06A	118701 118699				
-07A	2 118700				
-08A	3 118701				
-09A	118715 2				
-10A	6 3				
-11A	7 118715				
-12A	8 6				
-13A	Fm 11/24/10 7				
-14A	8				
-15A	9				

Comments:

Method: CARB430/Radiello 165Pressure: 137 bar

U s e	File #	Sample Name/Client	DF	Loaded by Init.	Date	Time	Reviewed by Init.	Comments
X	f1123001	CCV 1838-41-10ug ACU WASH ^{an112310}	1.00	Cal	11/23/10	0740	Cal	FT shift
✓	2	CCV 1838-41-10ug				0801		
✓	3	CCB				0822		
✓	4	Test Blank -1				0842		an112310 lot # 3 p.3 Notolue
✓	5	-2				0903		lot # p.99 ↓
✓	6	-3				0924		lot # p.3 w/ to luen
✓	7	↓ -4				0945		lot # p.99
✓	8	1838-41-10ug CCV				1006		
✓	9	CCB				1027		
✓	10	1011418A-Lab Blank				1048		
✓	11	1011418B-Lab Blank				1109		
X	12	1011418A-LCS				1130		low recoveries
X	13	↓ -LCSD				1150		↓
✓	14	1011418B-LCSD ^{an112310}				1211		
✓	15	↓ LCSD				1232		
✓	16	ACU WASH				1253		
✓	17	1838-41-10ug CCV				1314		
✓	18	CCB				1335		
✓	19	1011418A-LCS				1356		
✓	20	↓ -LCSD				1417		
✓	21	CARB430MDL-BLANK				1438		CARB430 extraction with Nucure 11/23/10
✓	22	-1				1458		spike 200ug Spike Std 1838-41-20ug/5ml
✓	23	-2				1519		
✓	24	-3				1540		
✓	25	-4				1601		
✓	26	-5				1622		spike 200ug Spike Std 1838-36-20ug/5ml
✓	27	-6				1643		
✓	28	-7				1704		

Calculation Check:

File ID: f1123023.d Compound: O-TolualdehydeInitials: Cal

Sample Amt = $\frac{\text{Area Counts Sample}}{\text{RF}} \times \text{DF} = \left(\frac{29082}{36466} \right) \times (1.00) =$

0.798

Reported Result:

0.798 0.798

an112410

Method: Carb430/Radiello/65Pressure: 137 bar

U s e	File #	Sample Name/Client	DF	Loaded by Init.	Date	Time	Reviewed by Init.	Comments
✓	H123029	CARB430MDL-8	1.00	Cee	11/23/10	1725	Cee	Spice 2000 al spike std 1838-36-20mg
✓	30	CCV1838-41-10mg				1746		
✓	31	CEB				1806		
C	32	1011298-06A				1827		C=confirmation Front impinger
C	33	↓ -07A				1848		back impinger
✓	34	1011299B-12A				1909		
✓	35	1011418A-05A				1930		
✓	36	↓ -0A				1951		
✓	37	↓ -15A				2012		
✓	38	1011418B-20A				2033		
✓	39	↓ -25A				2054		
✓	40	↓ -30A				2114		
✓	41	1011418A-04A				2135		
✓	42	CCV1838-41-10mg				2156		
✓	43	CEB				2217		
✓	44	1011418A-09A				2238		
✓	45	↓ -14A				2259		
✓	46	1011418B-19A				2320		
✓	47	↓ -24A			✓	2341		
✓	48	↓ -29A			11/24/10	0002		
✓	49	1011418A-01A	10.0			0022		
✓	50	↓ -02A				0043		
✓	51	↓ -03A				0104		
✓	52	↓ -06A				0125		
✓	53	↓ -07A				0146		
✓	54	1838-41-10mg CCV	1.00			0207		
✓	55	CEB				0228		
✓	↓ 56	1011418A-08A	10.0			0249		

Calculation Check:

File ID: _____ Compound: _____ Initials: _____

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

Cary Leaf

Signed

11/24/10

Date

Cee 11/24/10

Revised 07/08

Method: Carb430/Radiello 165Pressure: 137 bar

U s e	File #	Sample Name/Client	DF	Loaded by Init.	Date	Time	Reviewed by Init.	Comments
✓	f1123057	1011418A-11A	10.0	Cee	11/24/10	0310	Cee	
✓	58	↓ -12A	↓	↓		0330	↓	
✓	59	↓ -13A	↓	↓		0357	↓	
✓	60	1011418B-16A	↓	↓		0412	↓	
✓	61	↓ -17A	↓	↓		0433	↓	
✓	62	↓ -18A	↓	↓		0454	↓	
✓	63	↓ -21A	↓	↓		0515	↓	
✓	64	↓ -22A	↓	↓		0536	↓	
✓	65	↓ -23A	↓	↓		0557	↓	
✓	66	1838-41-10mg CCV	1.00	↓		0617	↓	
✓	67	CeB	↓	↓		0638	↓	
✓	68	1011418B-26A	10.0	↓		0659	↓	
✓	69	↓ -27A	↓	↓		0720	↓	
✓	70	↓ -28A	↓	↓		0741	↓	
✓	71	ACN WASH	1.00	↓		0802	↓	
✓	72	CCV 1838-41-10mg	↓	↓		0823	↓	

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:

Cary Leaf
Signed

11/24/10
Date

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: 11/11/10

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.
01A	118683	Air PASSIVE	Formaldehyde Analysis 11/1/10-11/5/10	13D 22H 10M
02A	118684			
03A	118685			
04A	118686			
05A	118687			
06A	118699		11/1/10-11/5/10	13D 23H 15M
07A	118700			
08A	118701			
09A	118702			
10A	118703			
11A	118715		11/1/10-11/5/10	13D 23H 20M
12A	118716			
13A	118717			
14A	118718			
15A	118719			

Special instructions:

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

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

Relinquished by: [Signature] of Environmental Health & Engineering, Inc. Date: 11/11/10
Received by: Aimee Wilkoff of (company name) Air Date: 11/18/10 09:00
Relinquished by: _____ of (company name) _____ Date: _____
Received by: _____ of (company name) _____ Date: _____
Relinquished by: _____ of (company name) _____ Date: _____
Received by: _____ of (company name) _____ Date: _____
Lab Data
Received by: _____ of Environmental Health & Engineering, Inc. Date: _____



FedEx

Page 1 of 2

SAMPLE RECEIPT SUMMARY

WORKORDER 1011418A

Client

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

Phone

800-825-5343

Fax

781-247-4305

Date Promised: 12/06/10

Date Completed: 12/6/10

Date Received: 11/18/10

PO#: 17131

Project#: 17131

Total \$: \$ 1,425.00

Sales Rep: TL

Logged By: MW

<u>Fraction</u>	<u>Sample #</u>	<u>Analysis</u>	<u>Collected</u>	<u>Amount\$</u>
01A	118683	Modified TO-11A	11/15/2010	\$90.00
02A	118684	Modified TO-11A	11/15/2010	\$90.00
03A	118685	Modified TO-11A	11/15/2010	\$90.00
04A	118686	Modified TO-11A	11/15/2010	\$90.00
05A	118687	Modified TO-11A	NA	\$90.00
06A	118699	Modified TO-11A	11/15/2010	\$90.00
07A	118700	Modified TO-11A	11/15/2010	\$90.00
08A	118701	Modified TO-11A	11/15/2010	\$90.00
09A	118702	Modified TO-11A	11/15/2010	\$90.00
10A	118703	Modified TO-11A	NA	\$90.00
11A	118715	Modified TO-11A	11/15/2010	\$90.00
12A	118716	Modified TO-11A	11/15/2010	\$90.00
13A	118717	Modified TO-11A	11/15/2010	\$90.00
14A	118718	Modified TO-11A	11/15/2010	\$90.00
15A	118719	Modified TO-11A	11/15/2010	\$90.00
16A	Lab Blank	Modified TO-11A	NA	\$0.00
17A	LCS	Modified TO-11A	NA	\$0.00
17AA	LCSD	Modified TO-11A	NA	\$0.00

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

SAMPLE RECEIPT SUMMARY Continued

Client

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

Phone

800-825-5343

Fax

781-247-4305

Date Promised:

Date Completed: 12/6/10

Date Received: 11/18/10

PO#: 17131

Project#: 17131

Total \$: \$ 1,425.00

Logged By: MW

Sales Rep:

<u>Fraction</u>	<u>Sample #</u>	<u>Analysis</u>	<u>Collected</u>	<u>Amount\$</u>
Misc. Charges	CVP (15) @ \$5.00 each.			\$75.00

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

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 #: 1011418A

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 <u>CIRCLE (YES / (NO))</u>
☐	☐	☒	☐	☐	☐	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) <u>10/11/9/10</u>
☐	☐	☒	☐	☐	☐	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:

LCS / LCSD

sig + sig/m3

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: <u>12/16/10</u>	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