

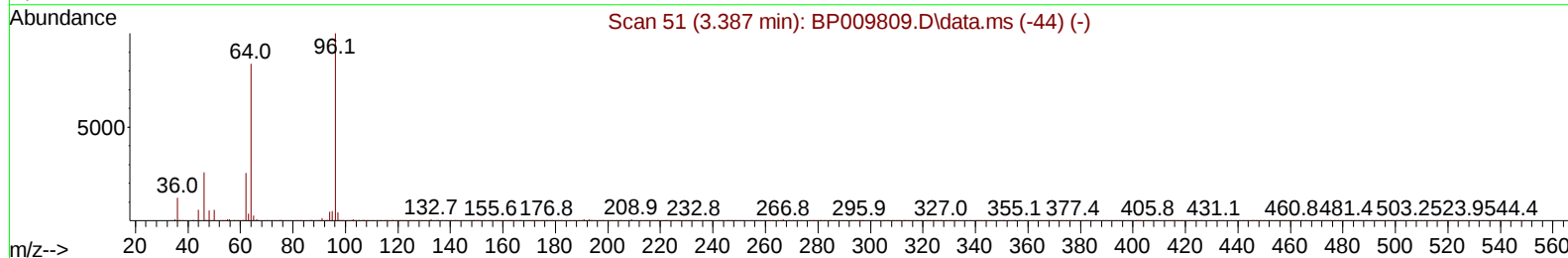
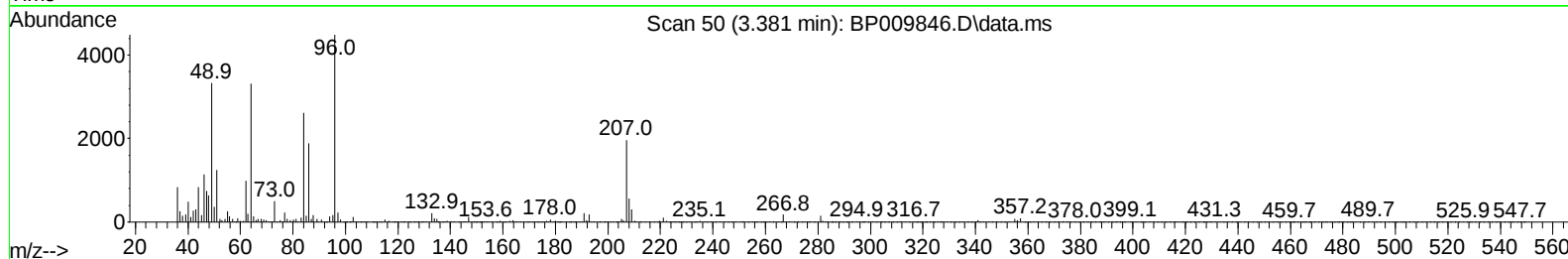
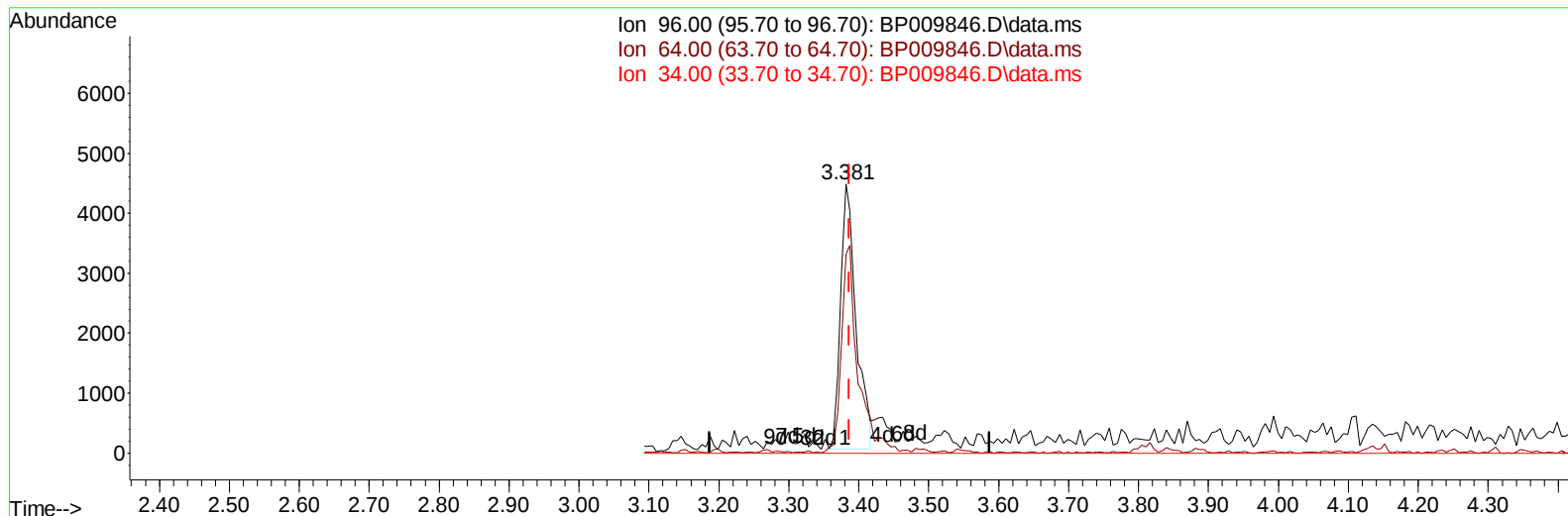
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009846.D
Acq On : 15 Apr 2022 01:58
Operator : CG/JU
Sample : N2293-14
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP2

Manual IntegrationsAPPROVED

Quant Time: Apr 15 04:21:52 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 13 02:15:22 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/15/2022
Supervised By :Yogesh Patel 04/22/2022



TIC: BP009846.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.381min (-0.006) 2.64 ng/uL

response 6907

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.88#
34.00	0.00	0.00
0.00	0.00	0.00

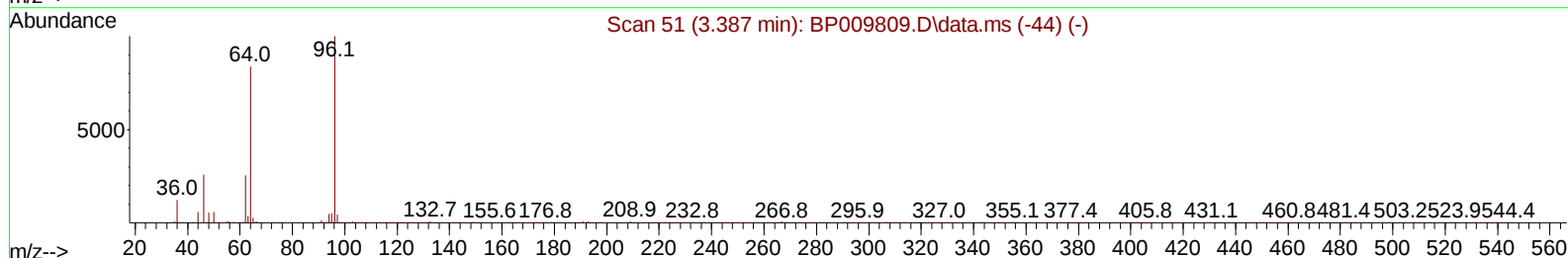
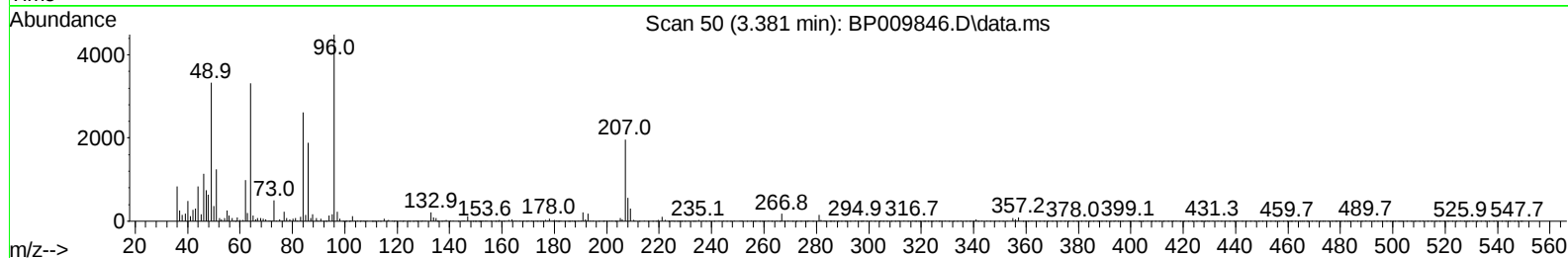
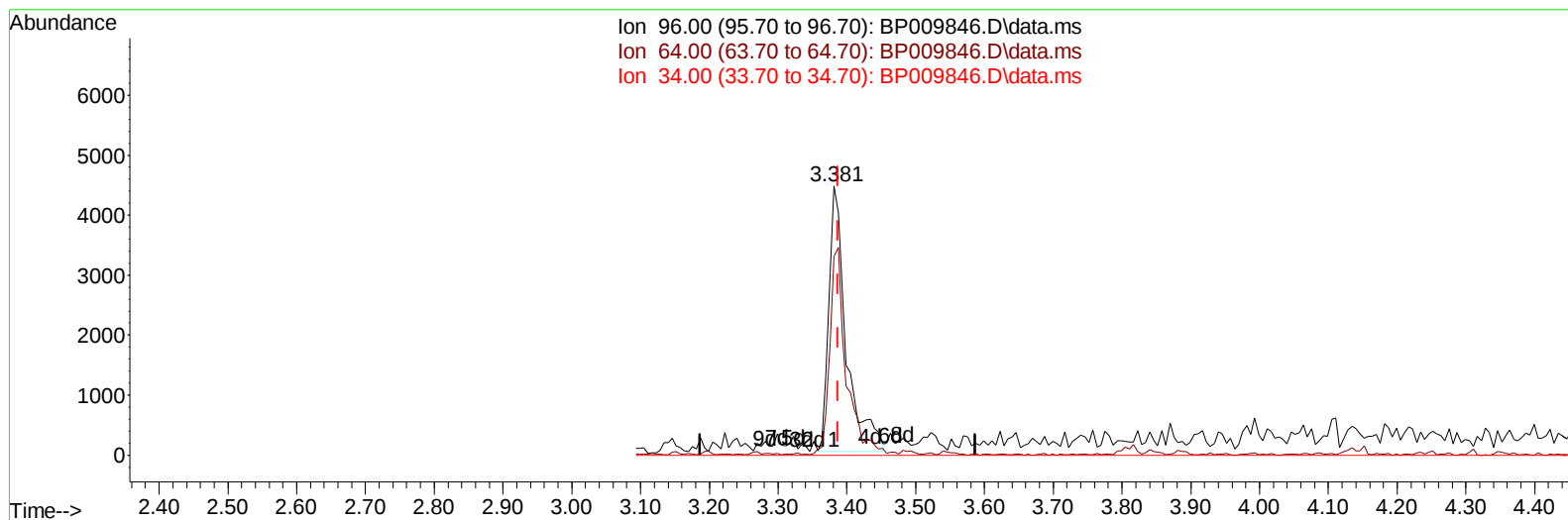
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TIC: BP009846.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.381min (-0.006) 2.97 ng/uL m

response 7777

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.88#
34.00	0.00	0.00
0.00	0.00	0.00

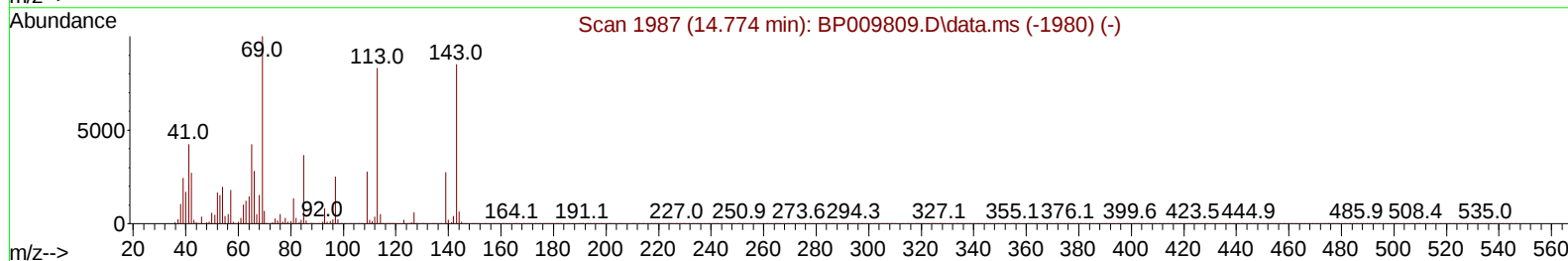
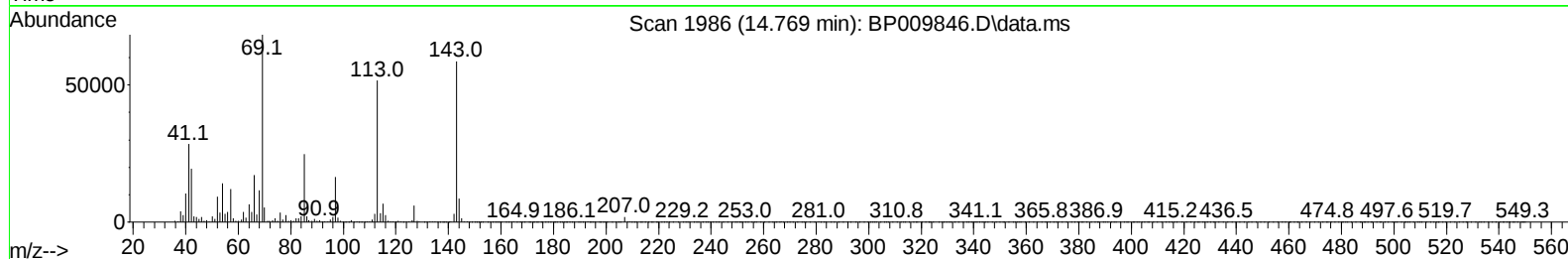
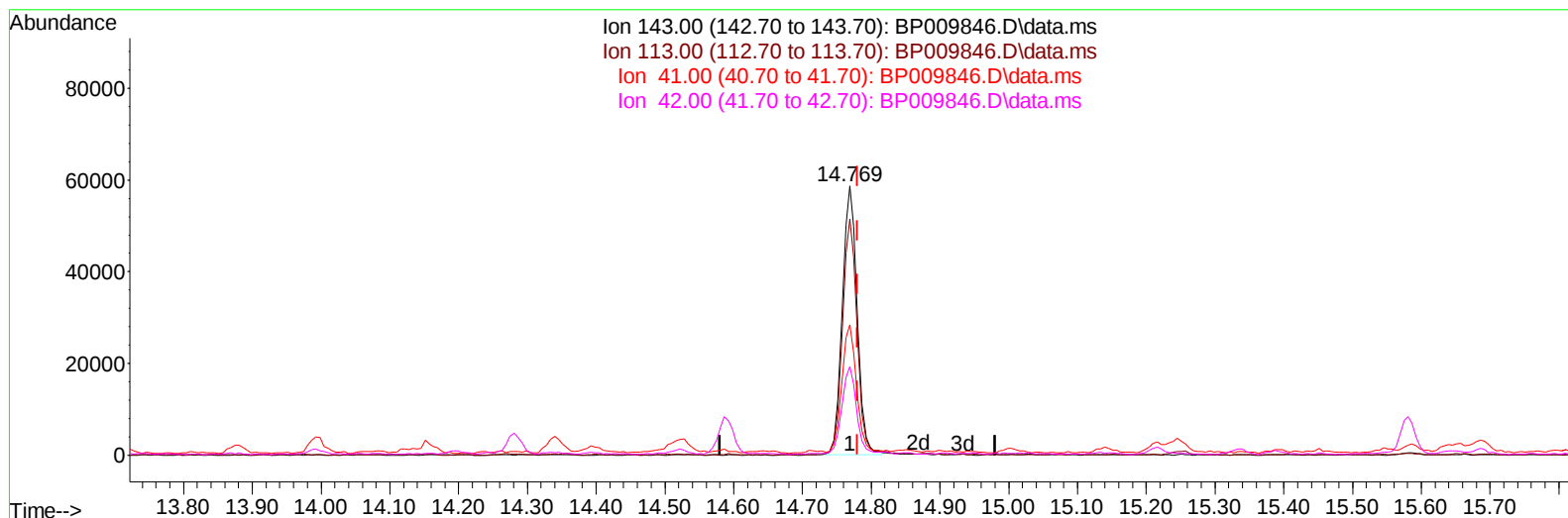
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TIC: BP009846.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.769min (-0.012) 16.67 ng/u1****response 87975**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	87.91#
41.00	23.20	48.48#
42.00	18.00	33.01#

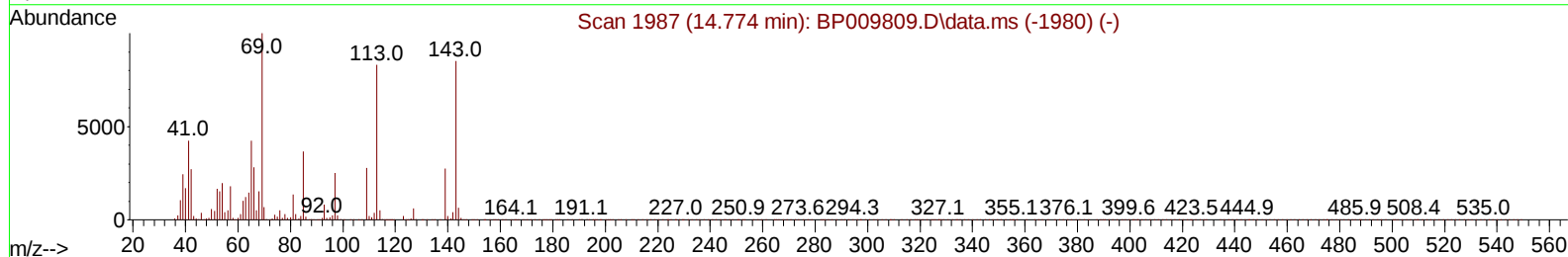
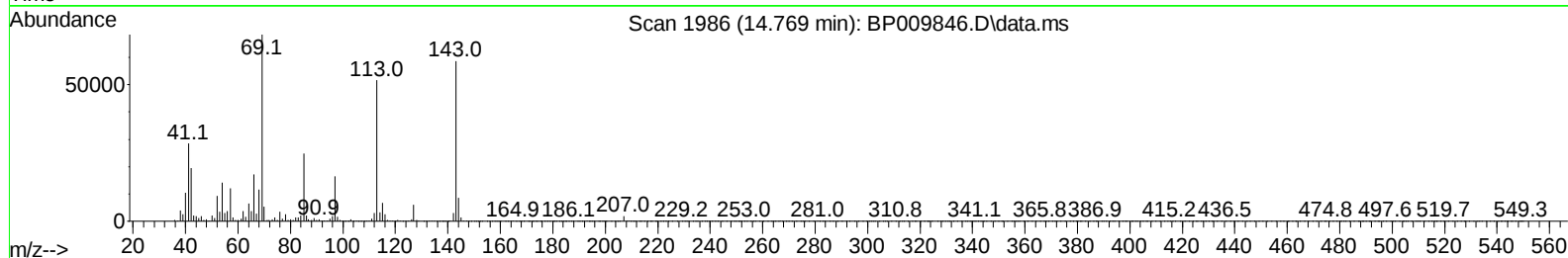
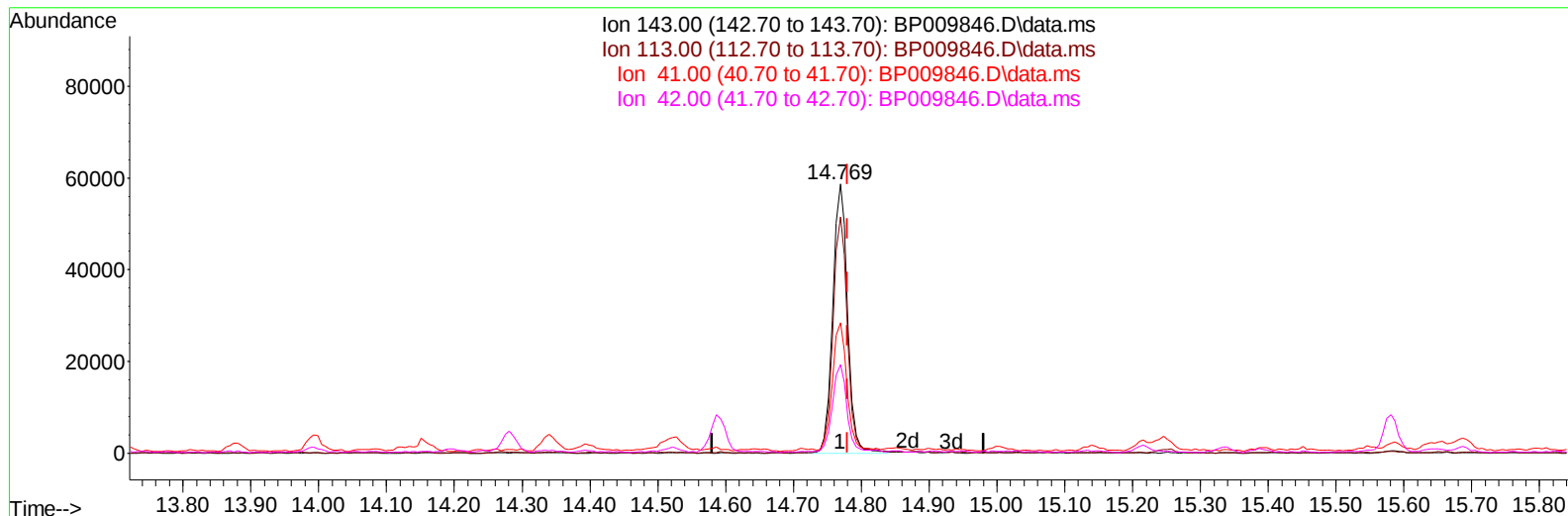
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 Operator : CG/JU
 Sample : N2293-14
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP009846.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (-0.012) 16.83 ng/ul m

response 88818

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	87.91#
41.00	23.20	48.48#
42.00	18.00	33.01#

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 Sample : N2293-14
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	116770	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.769	136	519116	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	370472	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	850072	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.416	240	954963	20.000	ng/ul	#-0.02
88) Perylene-d12	23.880	264	965122	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	7777m	2.972	ng/uL	0.00
4) Pyridine-d5	3.805	84	59141	8.198	ng/ul	0.00
7) Phenol-d5	7.111	99	150260	14.657	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.281	67	93644	15.344	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.487	132	113946	14.735	ng/ul	-0.01
15) 4-Methylphenol-d8	8.657	113	135088	15.785	ng/ul	-0.01
21) Nitrobenzene-d5	9.116	128	60848	16.254	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.840	143	63548	15.484	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.381	165	130027	15.140	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.893	131	167965	13.731	ng/ul	-0.01
46) Dimethylphthalate-d6	13.998	166	506092	17.391	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	513262	14.887	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.769	143	88818m	16.829	ng/ul	-0.01
60) Fluorene-d10	15.581	176	415812	16.985	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.687	200	44462	8.829	ng/ul	-0.01
73) Anthracene-d10	17.439	188	734713	18.052	ng/ul	-0.01
81) Pyrene-d10	19.663	212	959064	18.389	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.721	264	969460	19.289	ng/ul	-0.02
Target Compounds						
					Qvalue	
8) Phenol	7.140	94	21763	2.121	ng/ul#	90
18) 4-Methylphenol	8.716	108	18005	2.078	ng/ul	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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