

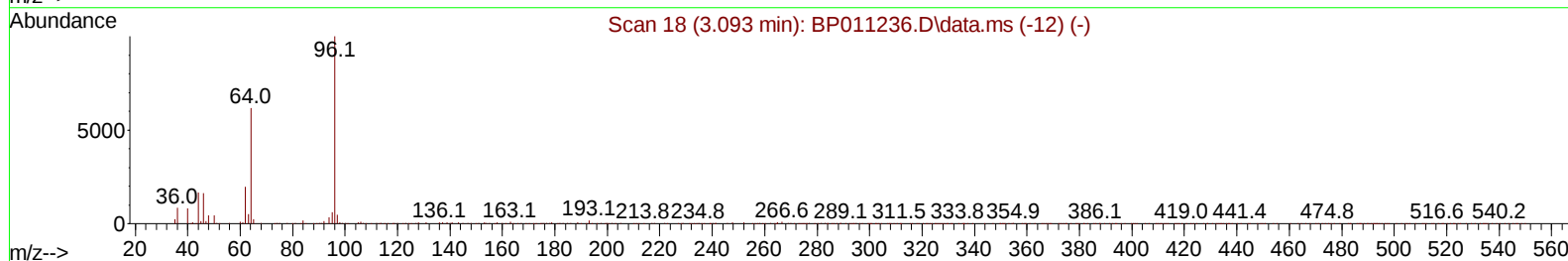
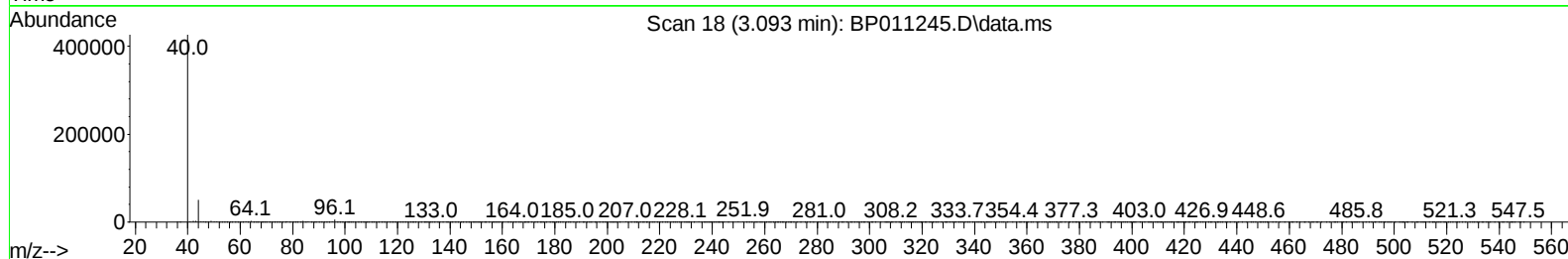
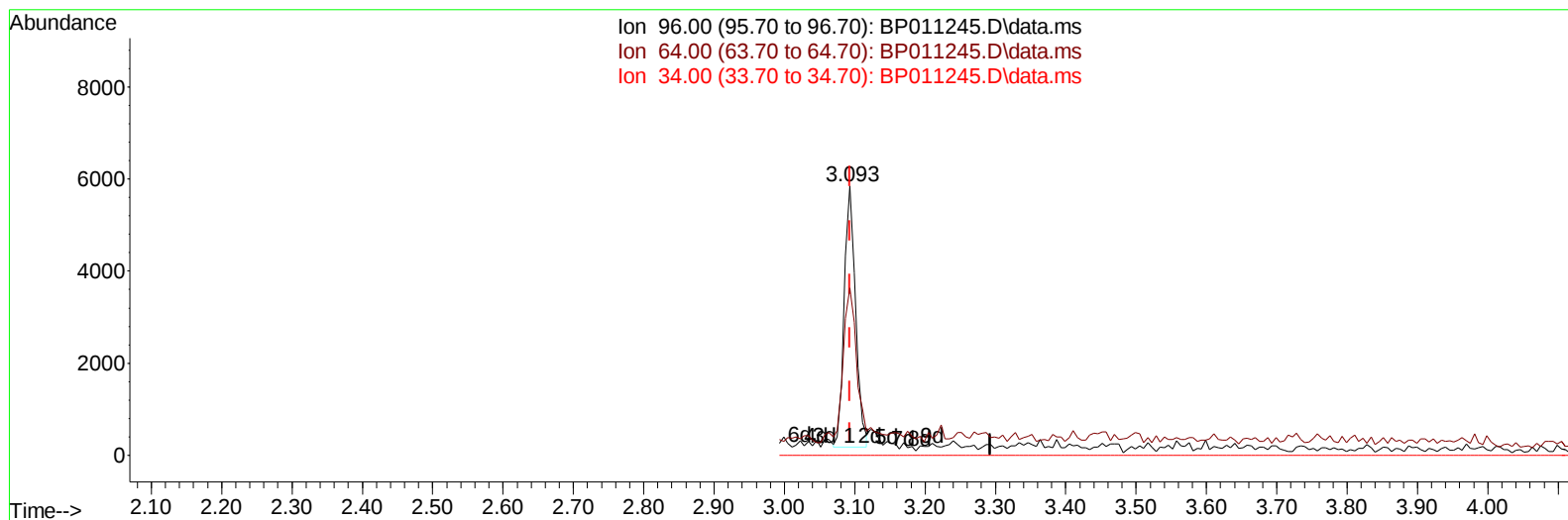
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080322\
Data File : BP011245.D
Acq On : 03 Aug 2022 15:13
Operator : CG/JU
Sample : N3900-16
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0CB3

Manual IntegrationsAPPROVED

Quant Time: Aug 03 23:00:17 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080222.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 02 23:52:37 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/04/2022
Supervised By :Sohil Jodhani 08/10/2022



TIC: BP011245.D\data.ms

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

3.093min (-0.000) 3.24 ng/uL

response 6275

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	62.13
34.00	0.00	0.00
0.00	0.00	0.00

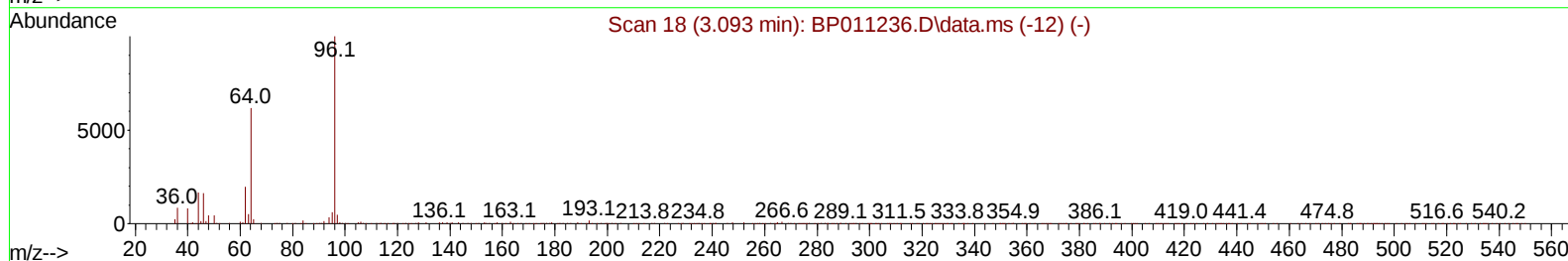
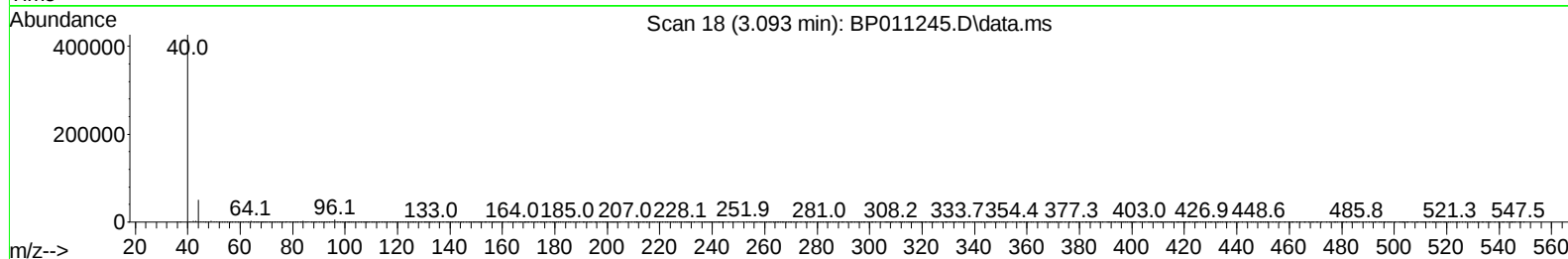
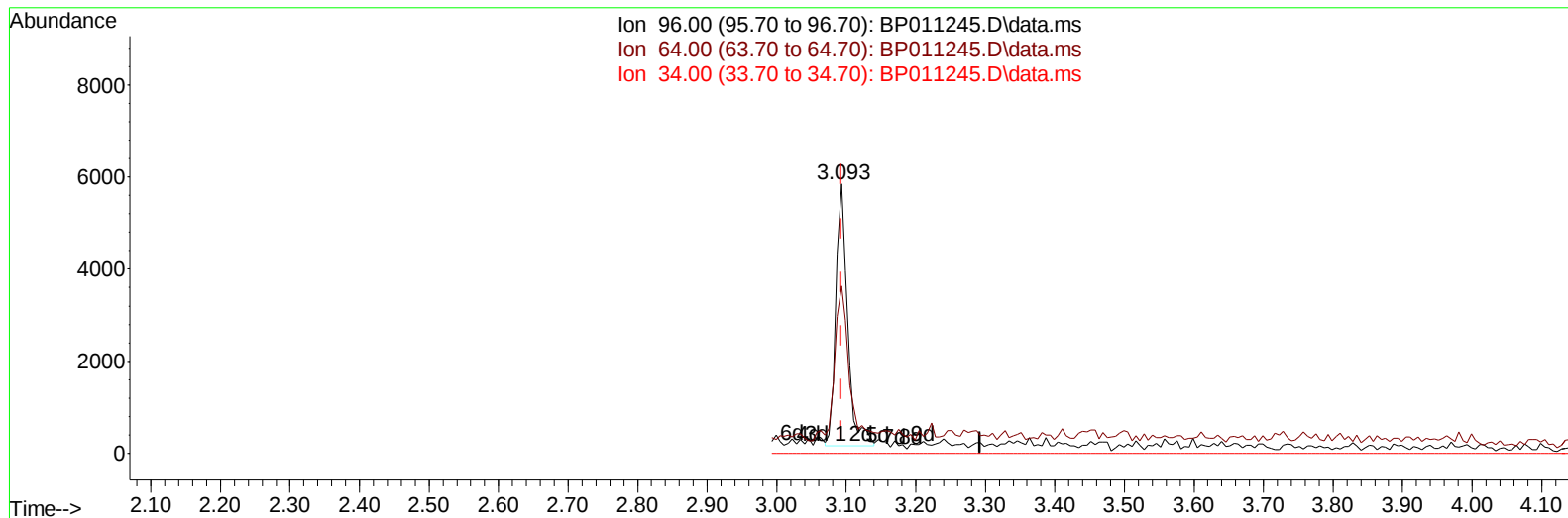
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TIC: BP011245.D\data.ms

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

3.093min (-0.000) 3.43 ng/uL m

response 6645

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	62.13
34.00	0.00	0.00
0.00	0.00	0.00

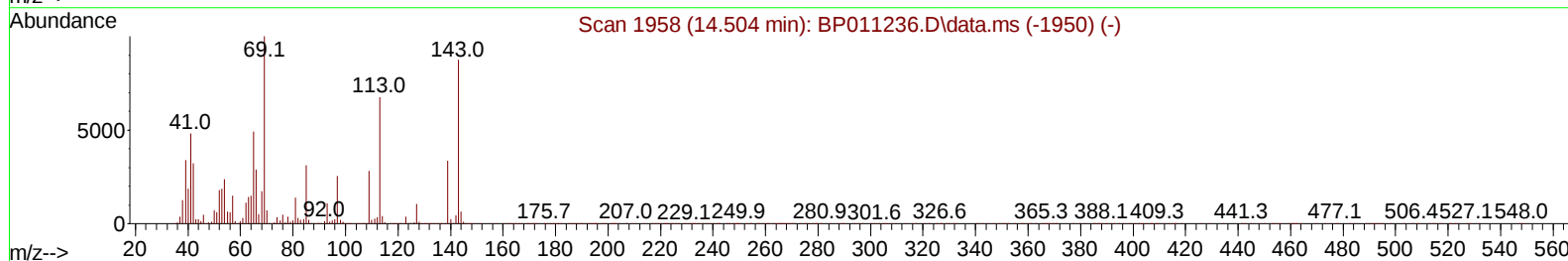
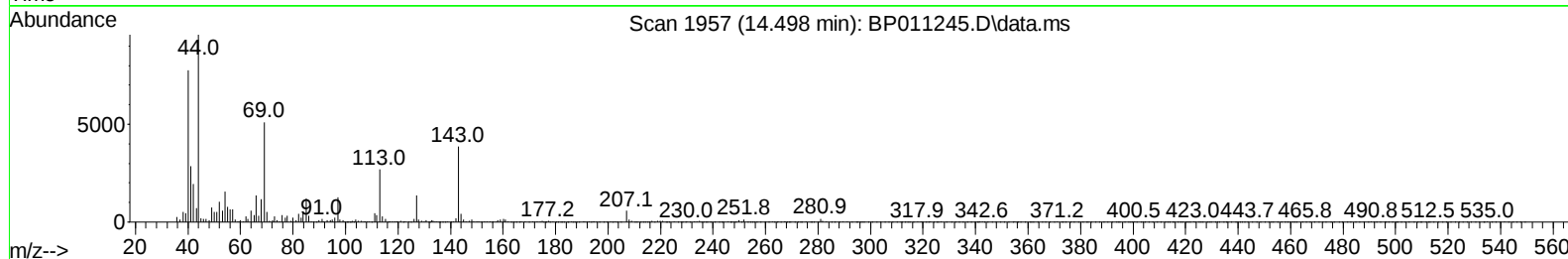
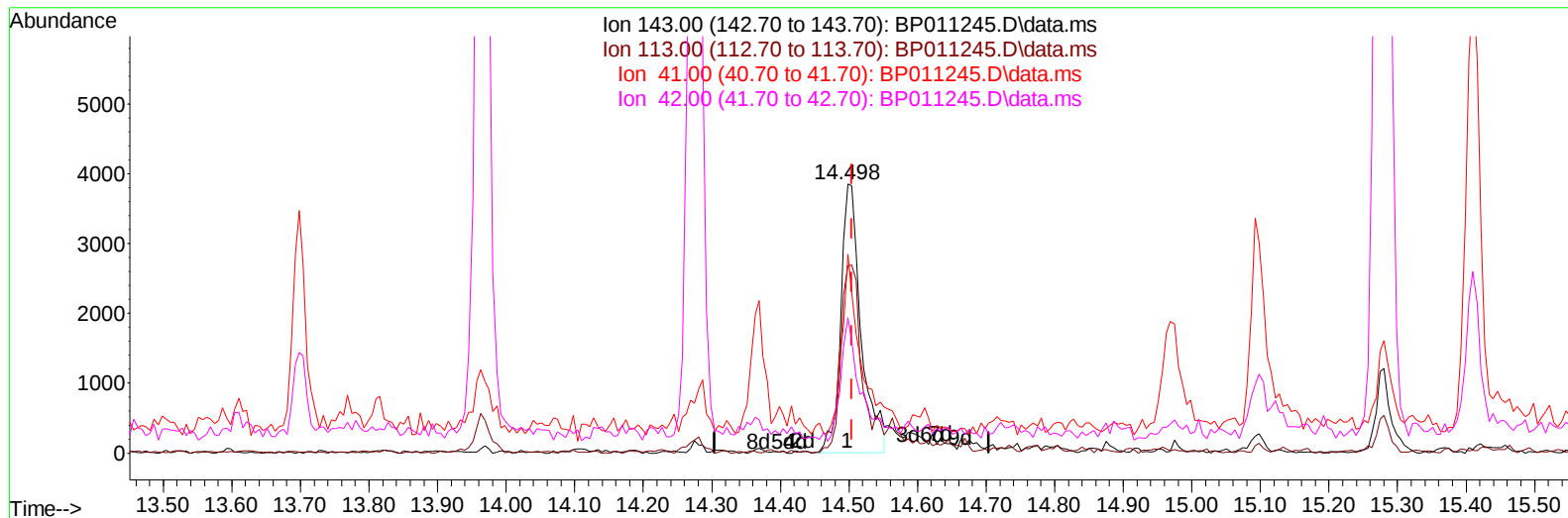
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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP011245.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.006) 1.84 ng/u1

response 7897

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	69.71
41.00	54.40	73.97#
42.00	35.50	50.42#

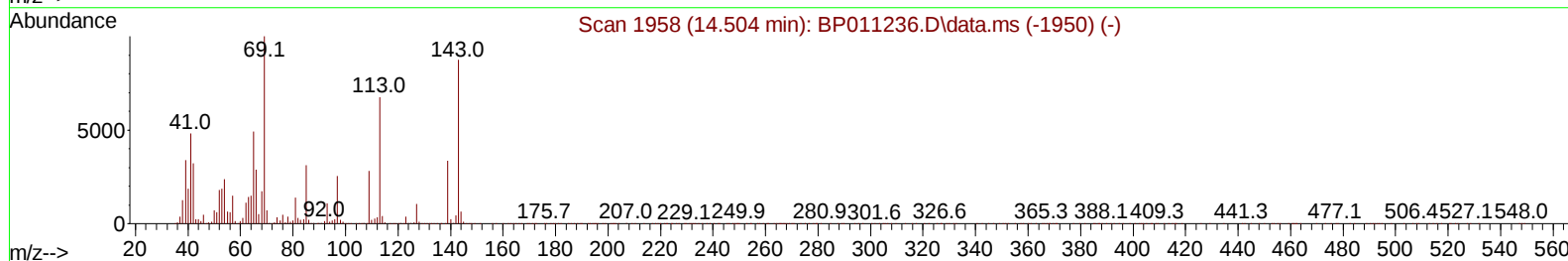
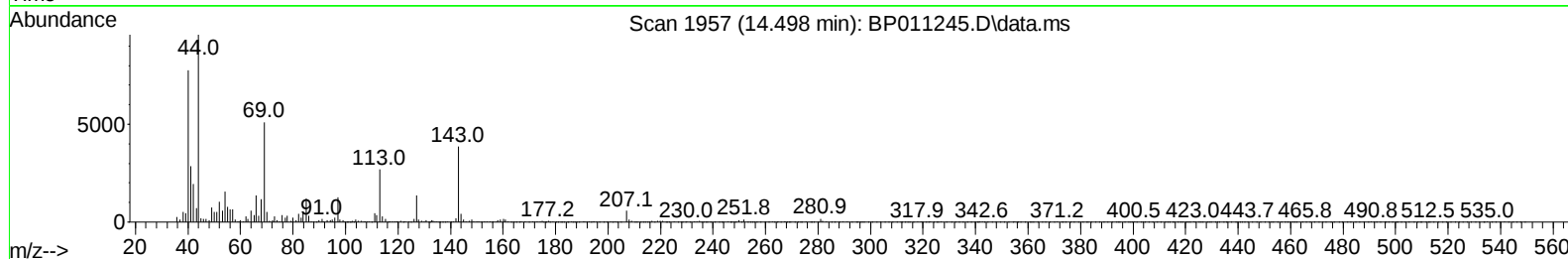
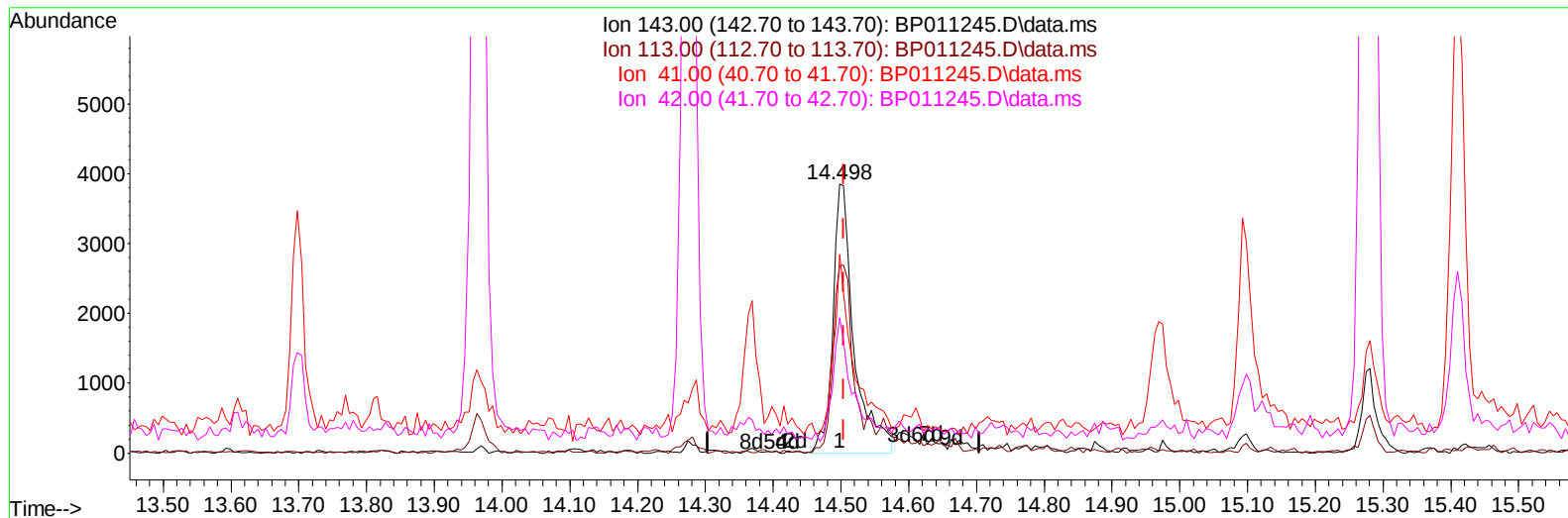
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 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
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Manual IntegrationsAPPROVED

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TIC: BP011245.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.006) 1.96 ng/ul m

response 8380

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	69.71
41.00	54.40	73.97#
42.00	35.50	50.42#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	156535	20.000	ng/ul	0.00
20) Naphthalene-d8	10.393	136	717043	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.275	164	367652	20.000	ng/ul	# 0.00
64) Phenanthrene-d10	17.045	188	667319	20.000	ng/ul	0.00
79) Chrysene-d12	21.174	240	710094	20.000	ng/ul	0.00
88) Perylene-d12	23.492	264	795836	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	6645m	3.432	ng/uL	0.00
4) Pyridine-d5	3.511	84	35949	6.884	ng/ul	0.00
7) Phenol-d5	6.781	99	54792	4.200	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	282283	29.988	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	190947	17.958	ng/ul	0.00
15) 4-Methylphenol-d8	8.322	113	111956	10.430	ng/ul	0.00
21) Nitrobenzene-d5	8.763	128	161205	31.256	ng/ul	0.00
24) 2-Nitrophenol-d4	9.487	143	106341	21.997	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	203260	22.048	ng/ul	0.00
31) 4-Chloroaniline-d4	10.546	131	329409	23.097	ng/ul	0.00
46) Dimethylphthalate-d6	13.698	166	859642	34.139	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	1105909	34.907	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	8380m	1.957	ng/ul	0.00
60) Fluorene-d10	15.281	176	810656	37.835	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	58558	18.898	ng/ul	0.00
73) Anthracene-d10	17.145	188	1225667	41.792	ng/ul	0.00
81) Pyrene-d10	19.410	212	1361762	37.929	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.345	264	1485316	37.453	ng/ul	0.00
Target Compounds						
						Qvalue
6) Benzaldehyde	6.752	77	23224	3.642	ng/ul	90
16) Acetophenone	8.434	105	26553	1.526	ng/ul	96

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

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