

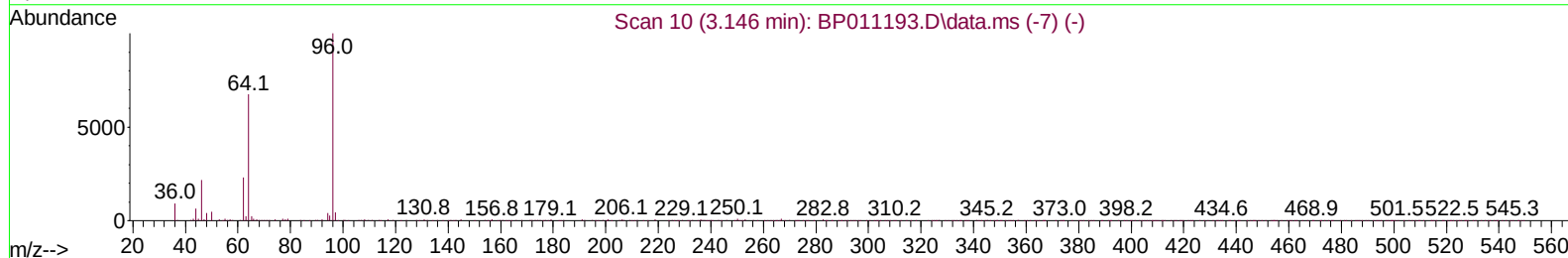
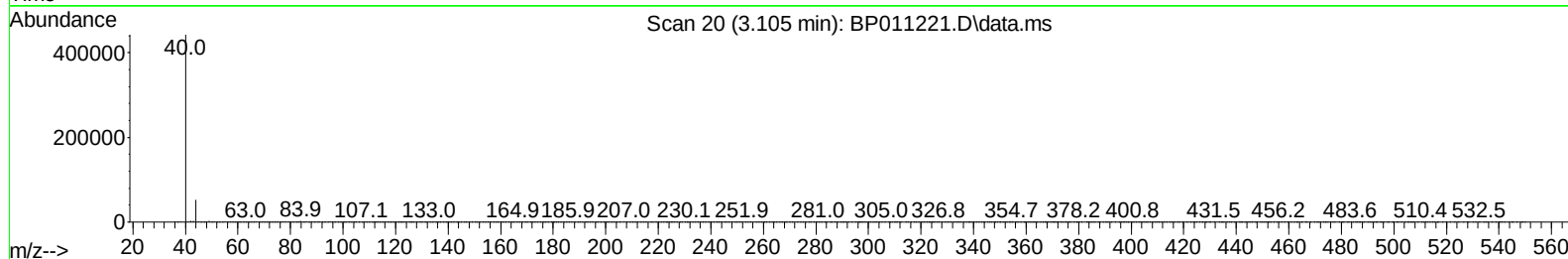
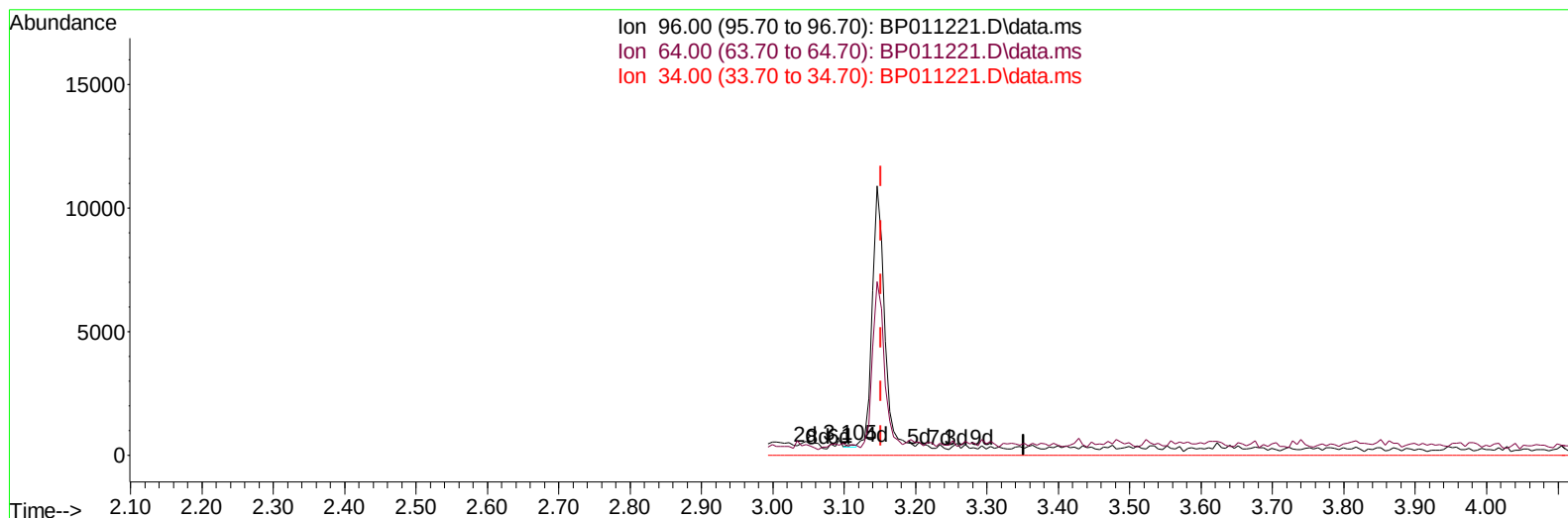
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080122\
 Data File : BP011221.D
 Acq On : 02 Aug 2022 08:57
 Operator : CG/JU
 Sample : N3900-10
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0C82

Manual IntegrationsAPPROVED

Quant Time: Aug 02 16:04:20 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 28 16:05:22 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/02/2022
 Supervised By :mohammad ahmed 08/03/2022



TIC: BP011221.D\data.ms

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

3.105min (-0.047) 0.02 ng/uL

response 100

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	84.12
34.00	0.00	0.00
0.00	0.00	0.00

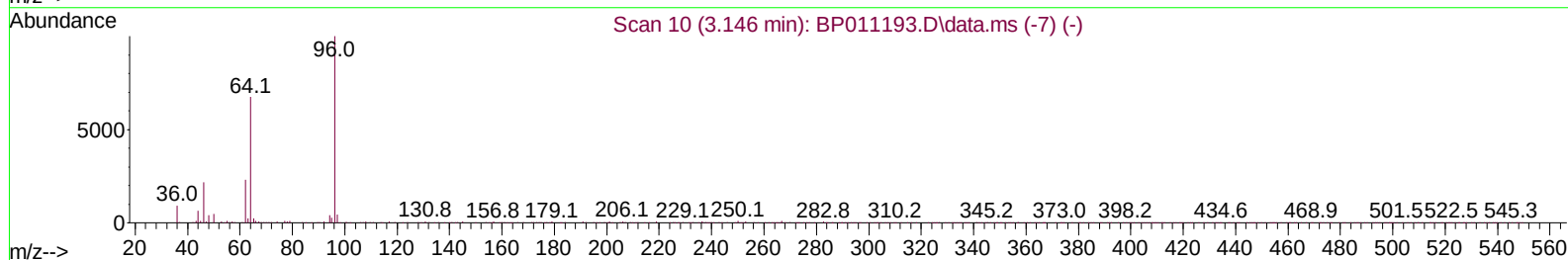
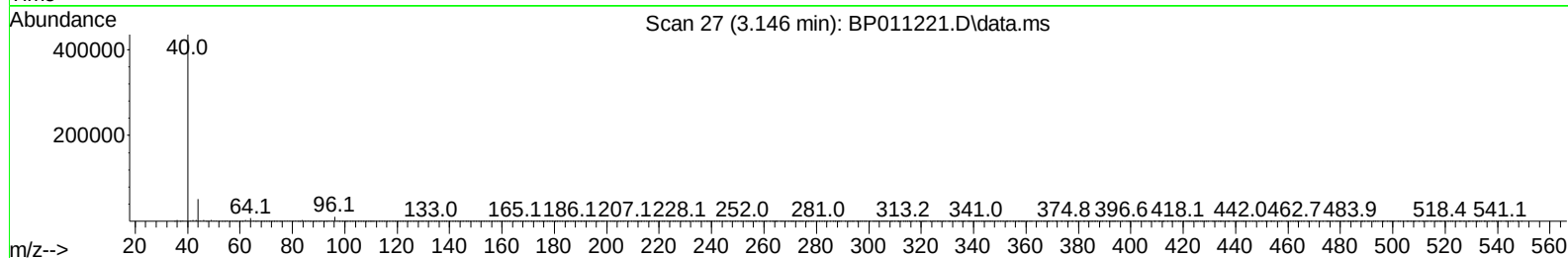
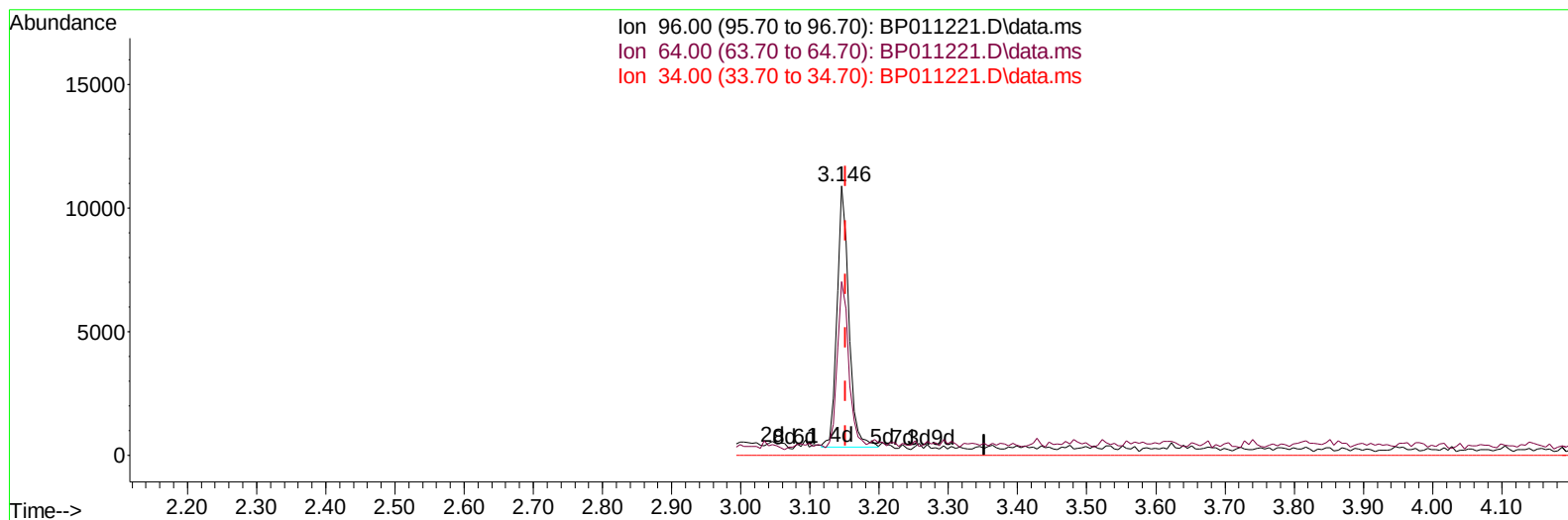
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TIC: BP011221.D\data.ms

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

3.146min (-0.006) 2.42 ng/uL m

response 12414

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	64.53#
34.00	0.00	0.00
0.00	0.00	0.00

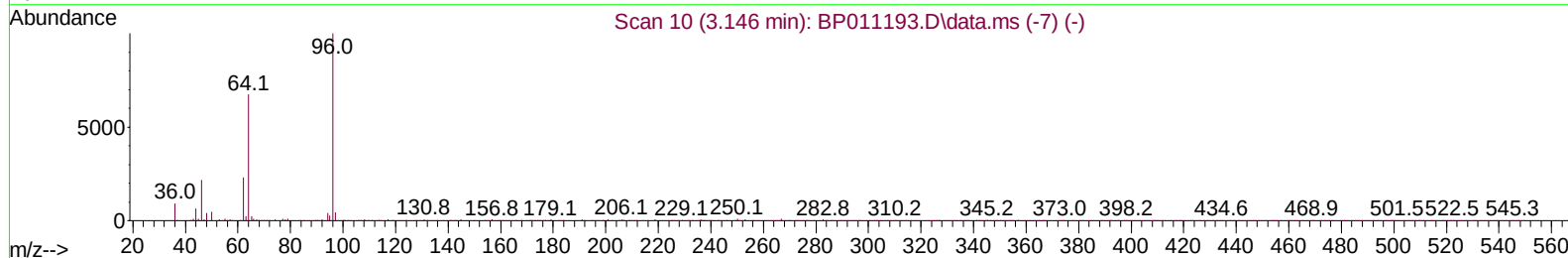
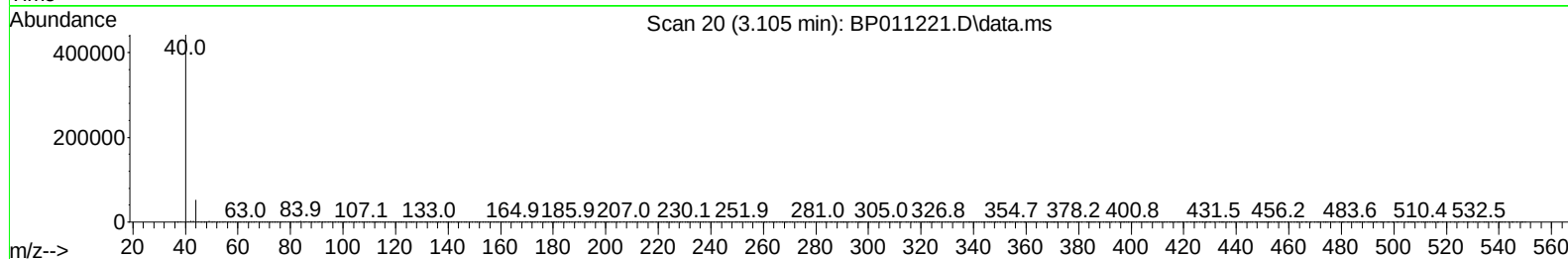
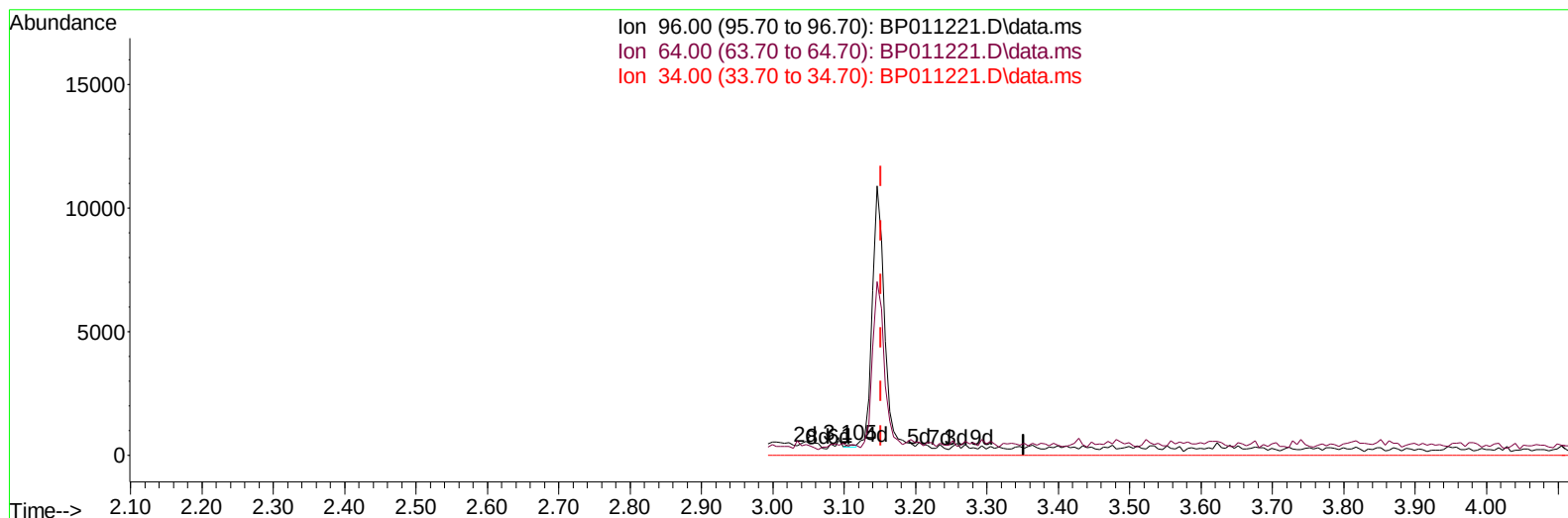
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 Operator : CG/JU
 Sample : N3900-10
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 C0C82

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TIC: BP011221.D\data.ms

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

3.105min (-0.047) 0.02 ng/uL

response 100

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	84.12
34.00	0.00	0.00
0.00	0.00	0.00

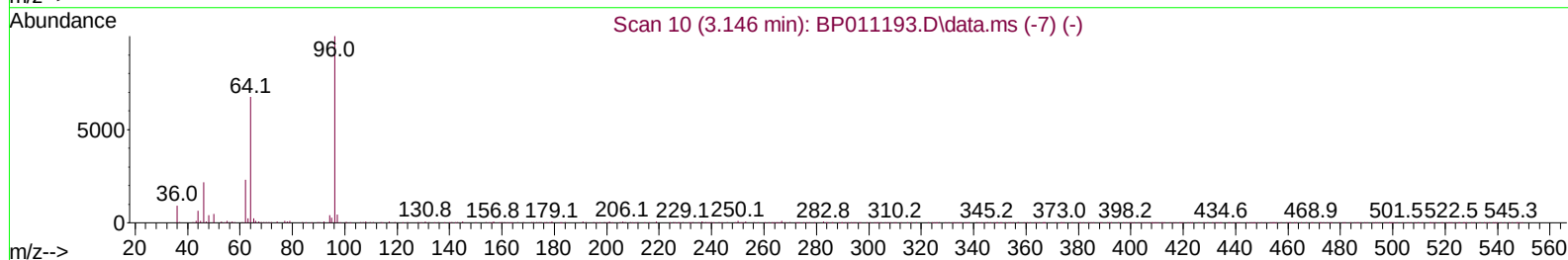
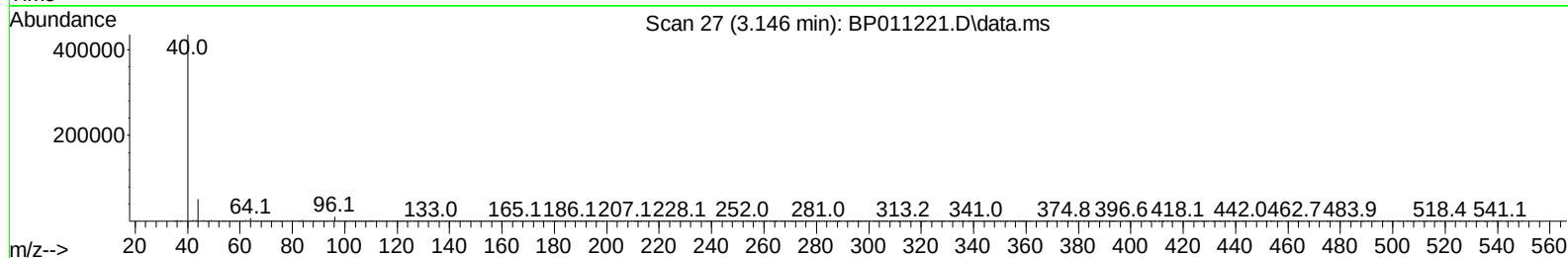
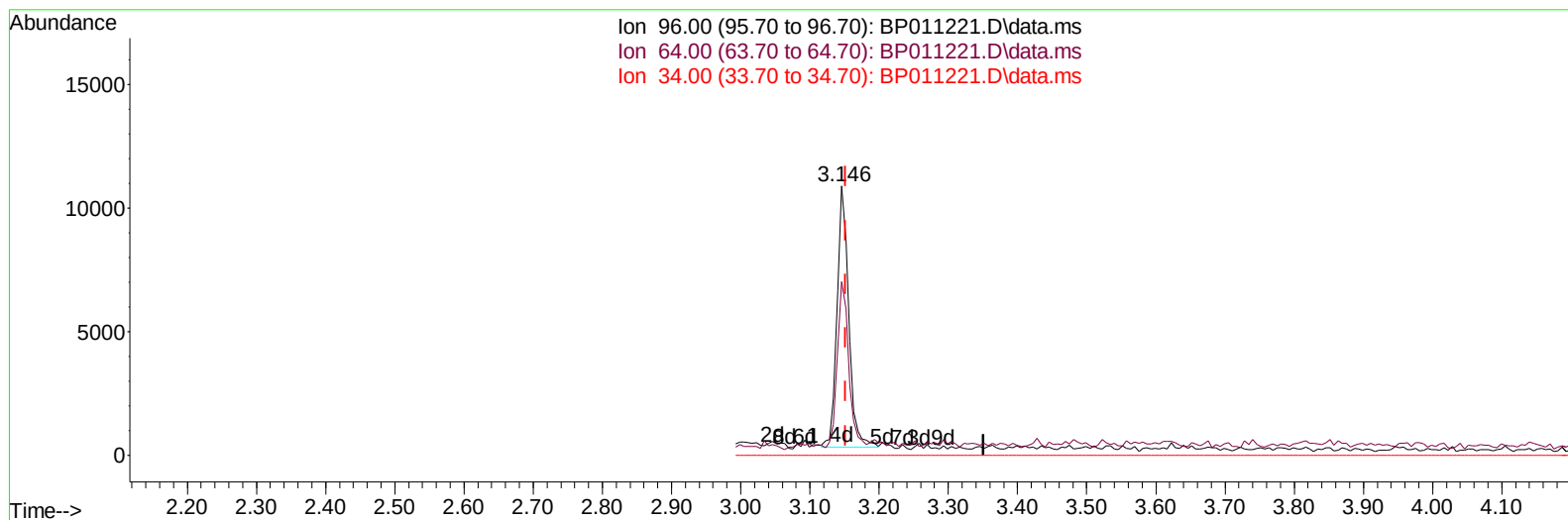
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 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
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TIC: BP011221.D\data.ms

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

3.146min (-0.006) 2.41 ng/uL m

response 12350

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	64.53#
34.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.646	152	234479	20.000	ng/ul	0.00
20) Naphthalene-d8	10.440	136	981789	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.316	164	496756	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.086	188	955216	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.215	240	946782	20.000	ng/ul	# 0.00
88) Perylene-d12	23.562	264	1013113	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	12350m	2.412	ng/uL	0.00
4) Pyridine-d5	3.563	84	54658	3.773	ng/ul	0.00
7) Phenol-d5	6.828	99	84115	4.110	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	317111	21.912	ng/ul	0.00
11) 2-Chlorophenol-d4	7.175	132	254271	15.765	ng/ul	0.00
15) 4-Methylphenol-d8	8.363	113	149978	8.958	ng/ul	0.00
21) Nitrobenzene-d5	8.810	128	173361	22.721	ng/ul	0.00
24) 2-Nitrophenol-d4	9.528	143	141761	19.396	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.063	165	235081	18.251	ng/ul	0.00
31) 4-Chloroaniline-d4	10.587	131	365549	17.507	ng/ul	0.00
46) Dimethylphthalate-d6	13.734	166	943712	27.884	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	1119920	26.332	ng/ul	0.00
54) 4-Nitrophenol-d4	14.539	143	10709	1.759	ng/ul	0.00
60) Fluorene-d10	15.316	176	788096	27.711	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	68198	14.163	ng/ul	0.00
73) Anthracene-d10	17.186	188	1366680	32.471	ng/ul	0.00
81) Pyrene-d10	19.445	212	1574645	30.951	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.410	264	1637388	32.539	ng/ul	0.00

Target Compounds Qvalue

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

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