

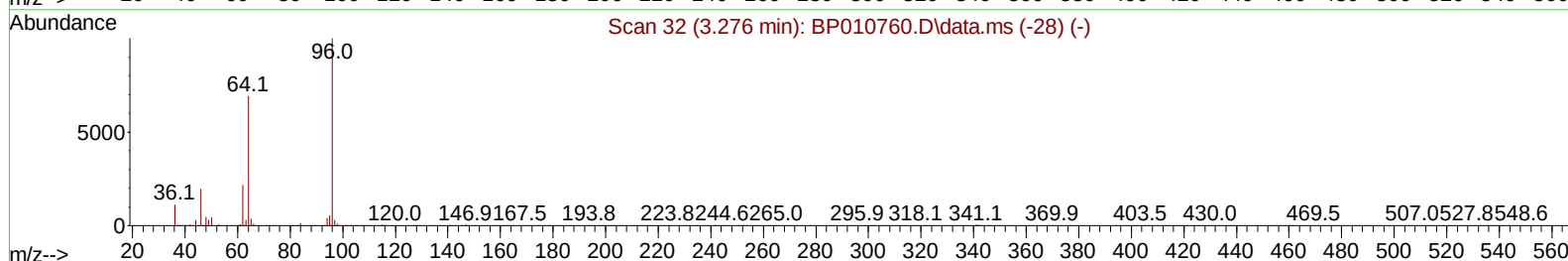
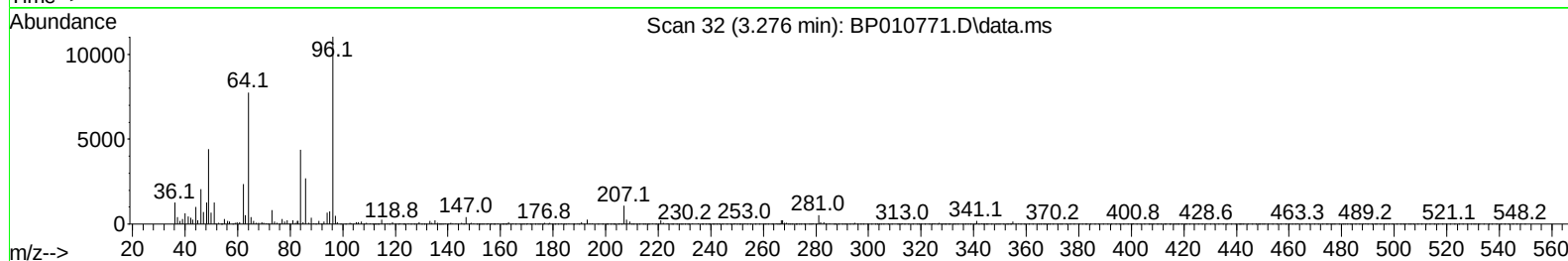
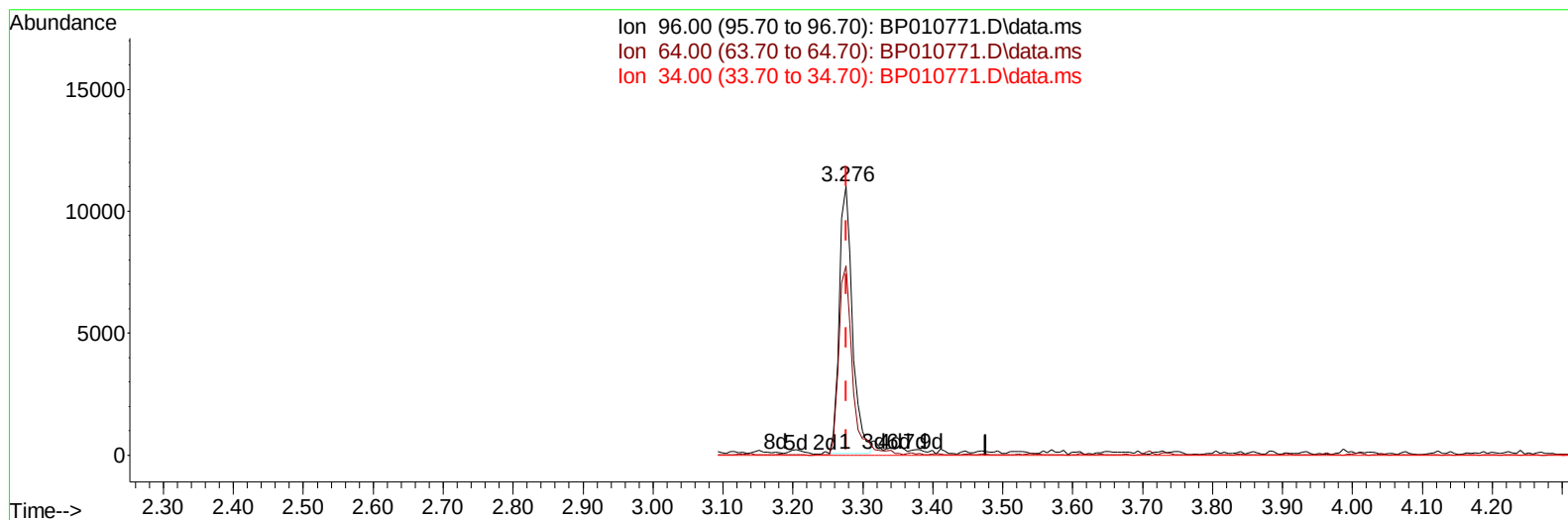
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062022\
 Data File : BP010771.D
 Acq On : 20 Jun 2022 20:47
 Operator : CG/JU
 Sample : N3358-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW4R6

Manual IntegrationsAPPROVED

Quant Time: Jun 21 00:47:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 20 16:18:42 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2022
 Supervised By :mohammad ahmed 06/27/2022



TIC: BP010771.D\data.ms

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

3.276min (-0.000) 4.95 ng/uL

response 14491

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	70.28
34.00	0.00	0.00
0.00	0.00	0.00

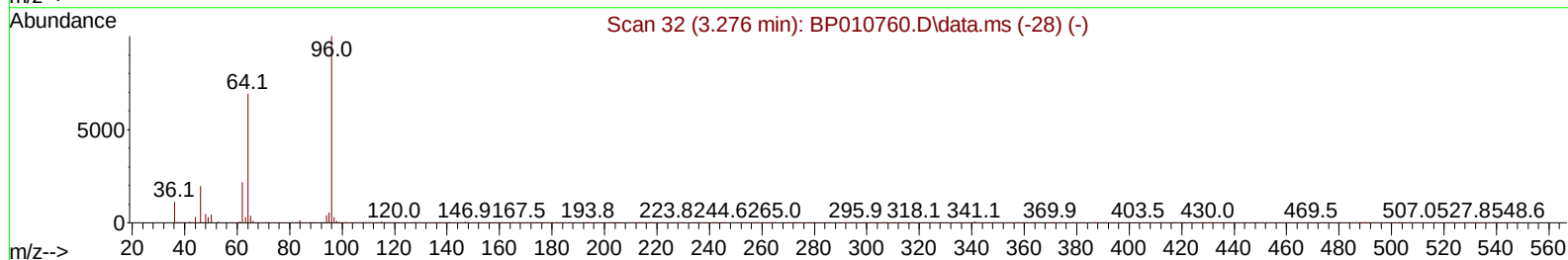
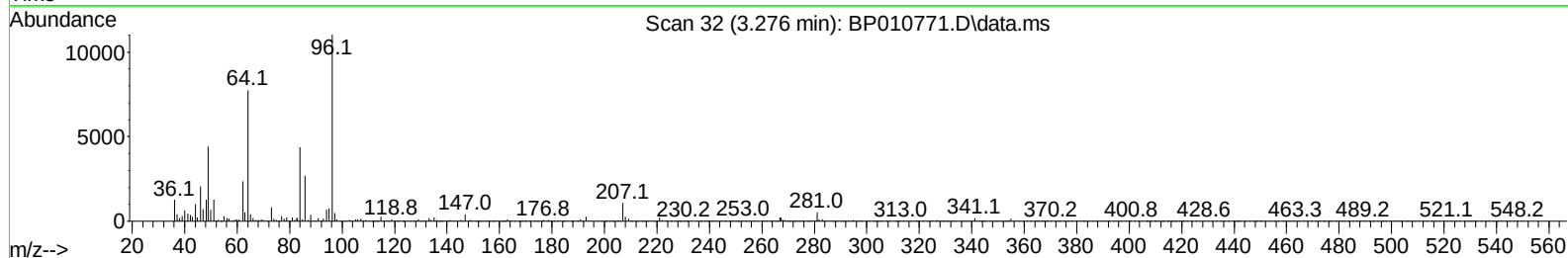
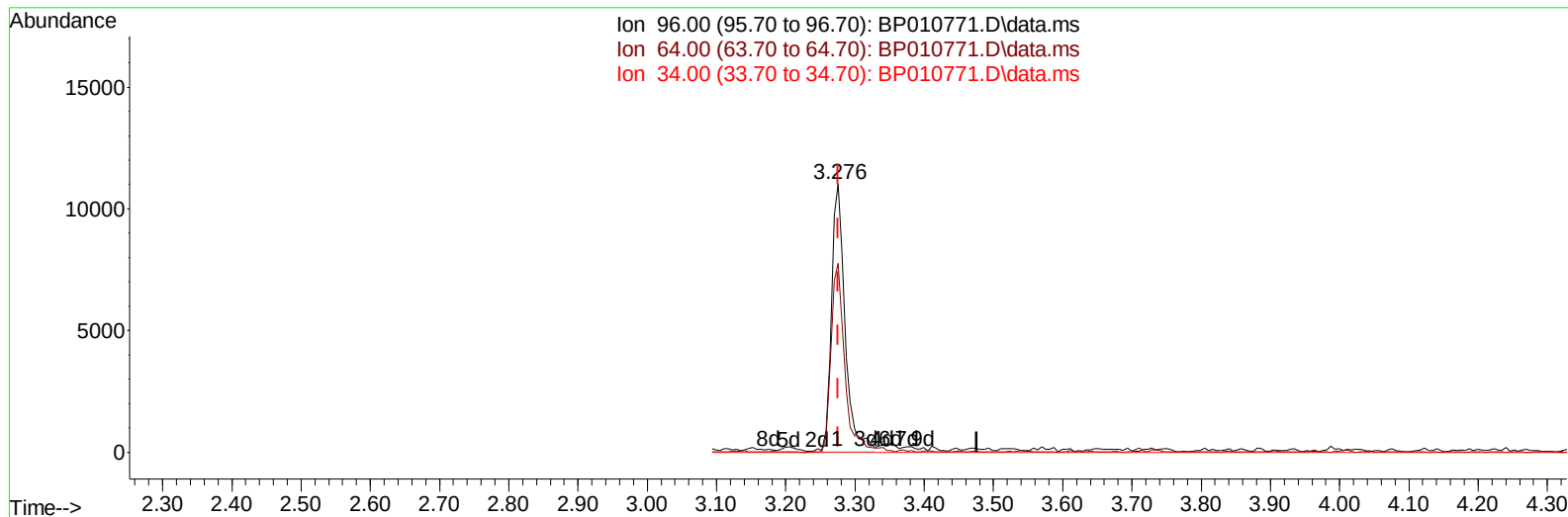
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TIC: BP010771.D\data.ms

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

3.276min (-0.000) 5.11 ng/uL m

response 14971

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	70.28
34.00	0.00	0.00
0.00	0.00	0.00

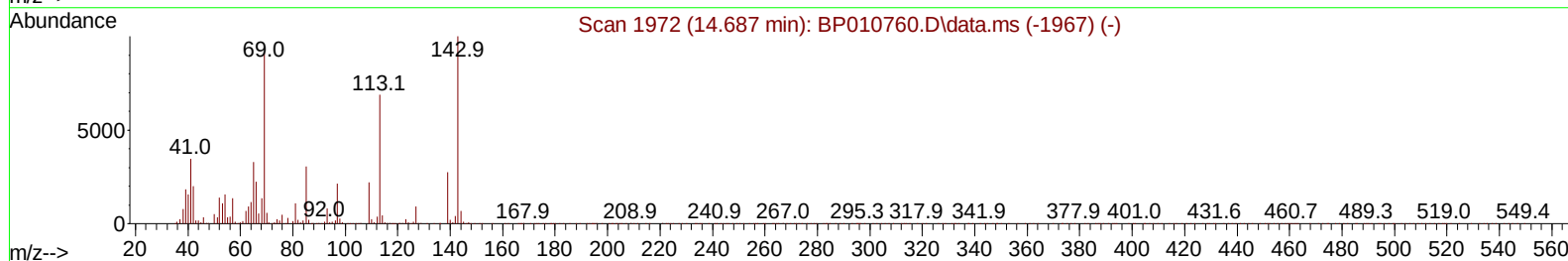
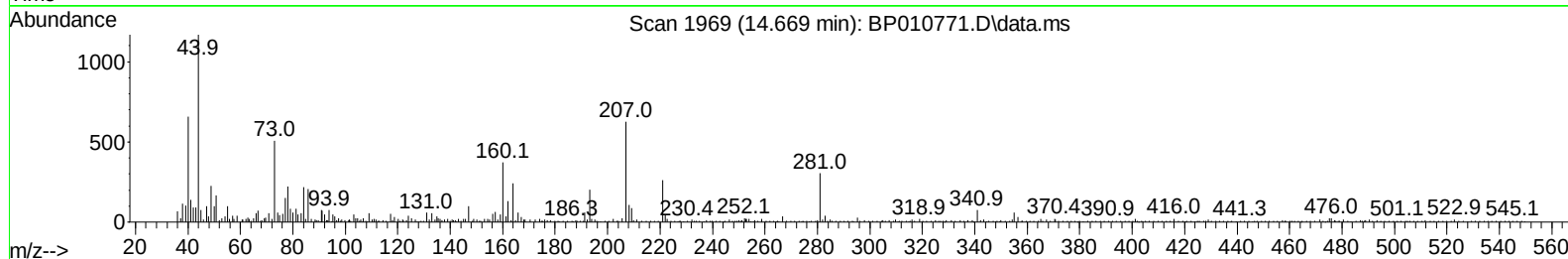
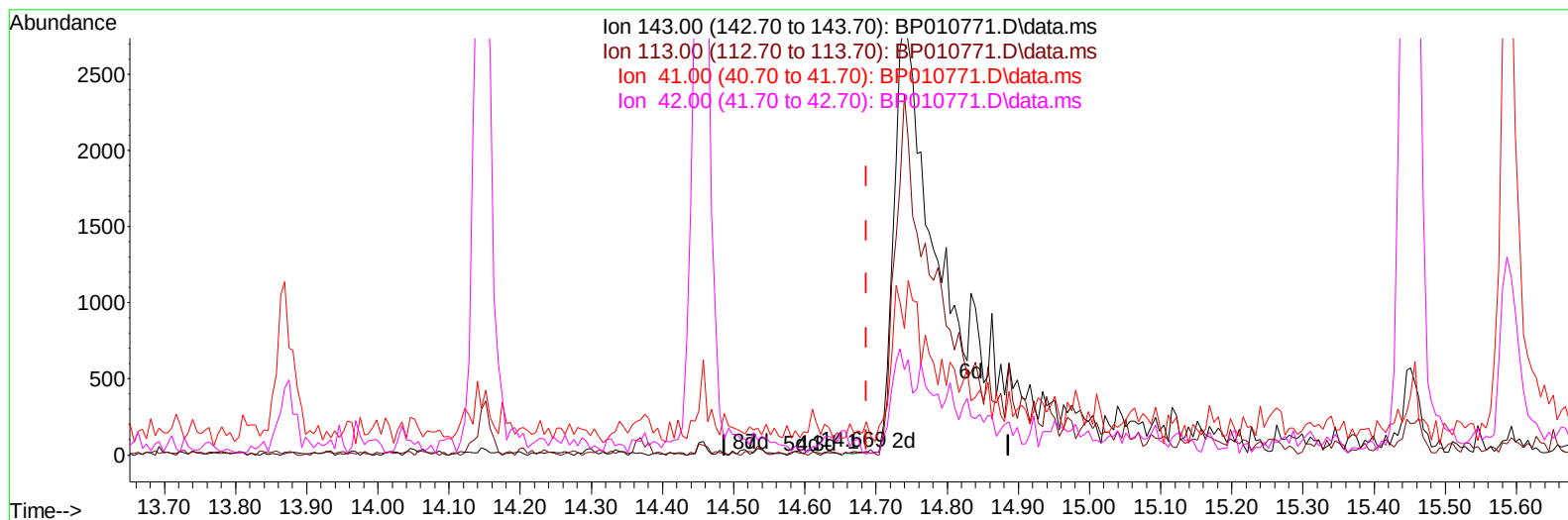
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TIC: BP010771.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.018) 0.00 ng/u1

response 11

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	47.62#
41.00	59.00	661.90#
42.00	41.10	442.86#

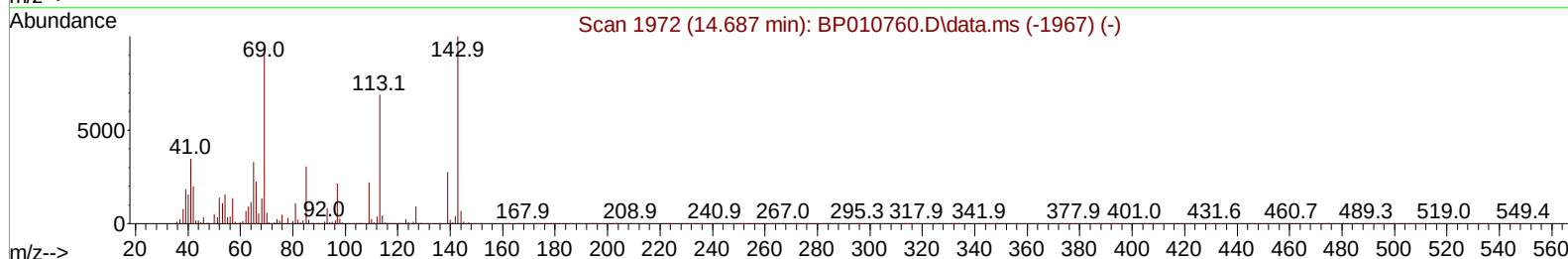
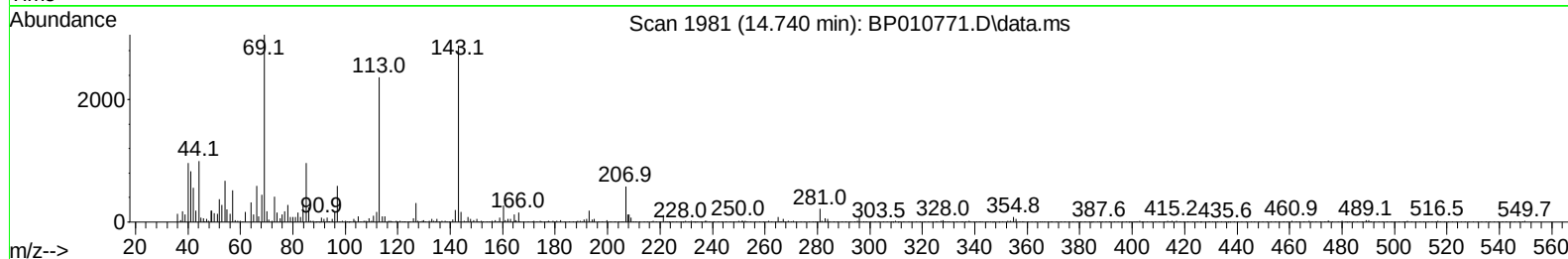
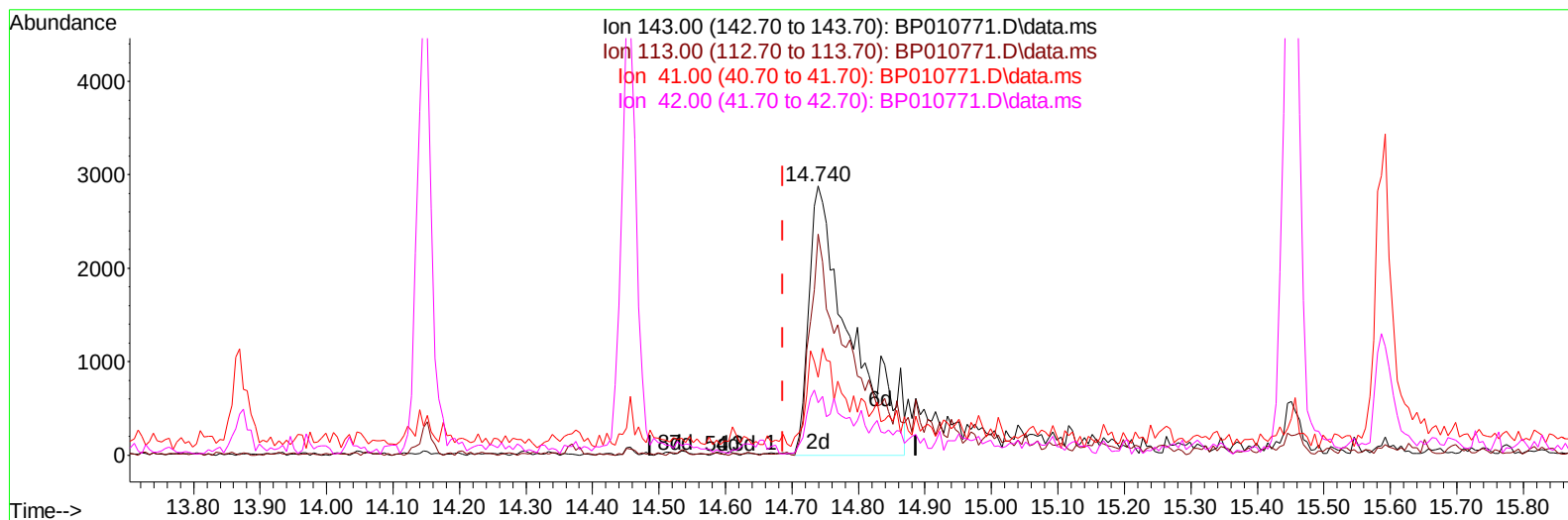
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TIC: BP010771.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.053) 4.33 ng/ul m

response 12658

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	82.08
41.00	59.00	28.83#
42.00	41.10	19.49#

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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	7.816	152	125891	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	463943	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	288104	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.210	188	590438	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	653250	20.000	ng/ul	0.00
88) Perylene-d12	23.692	264	735404	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	14971m	5.111	ng/uL	0.00
4) Pyridine-d5	3.705	84	66152	8.684	ng/ul	0.01
7) Phenol-d5	6.999	99	49136	5.379	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	148123	26.632	ng/ul	0.00
11) 2-Chlorophenol-d4	7.352	132	165347	21.267	ng/ul	0.00
15) 4-Methylphenol-d8	8.540	113	97286	13.181	ng/ul	0.00
21) Nitrobenzene-d5	8.975	128	98390	28.563	ng/ul	0.00
24) 2-Nitrophenol-d4	9.699	143	101321	27.455	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.252	165	157421	23.022	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	222343	23.084	ng/ul	0.00
46) Dimethylphthalate-d6	13.869	166	623328	31.513	ng/ul	0.00
49) Acenaphthylene-d8	14.146	160	694831	29.006	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	12658m	4.332	ng/ul	0.05
60) Fluorene-d10	15.451	176	550299	31.298	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.587	200	86522	24.561	ng/ul	0.00
73) Anthracene-d10	17.310	188	926194	34.590	ng/ul	0.00
81) Pyrene-d10	19.551	212	1194872	36.794	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.539	264	1356762	37.051	ng/ul	0.00

Target Compounds Qvalue

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

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