

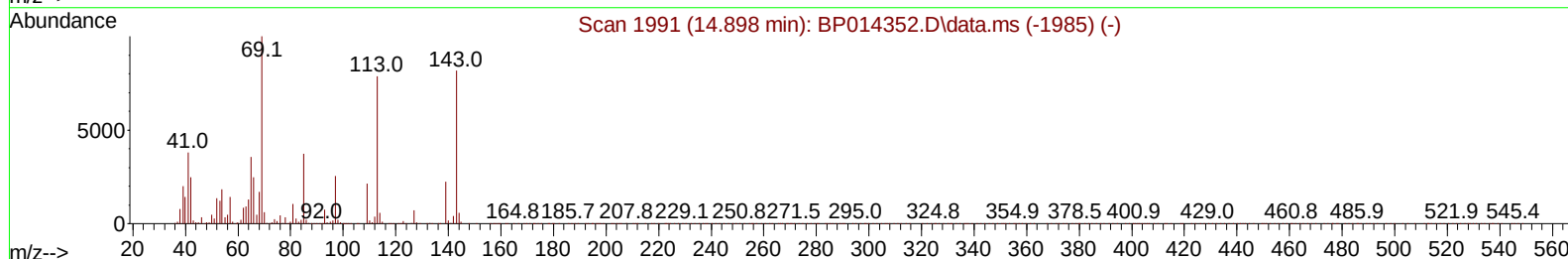
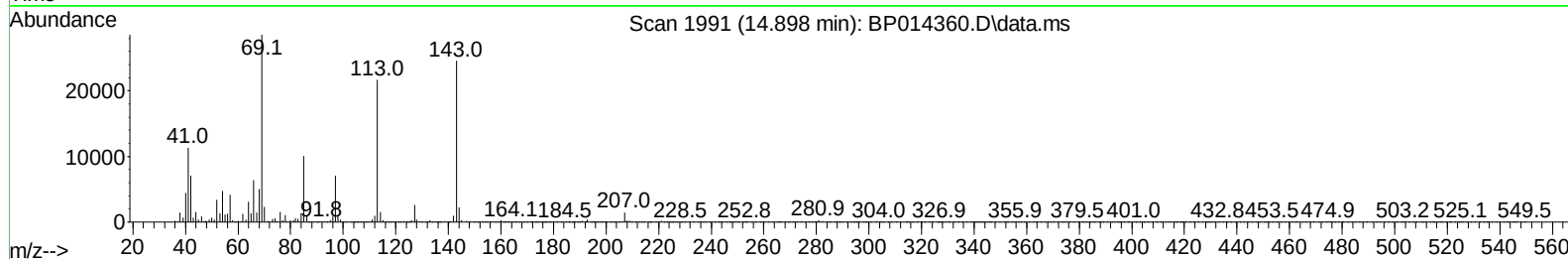
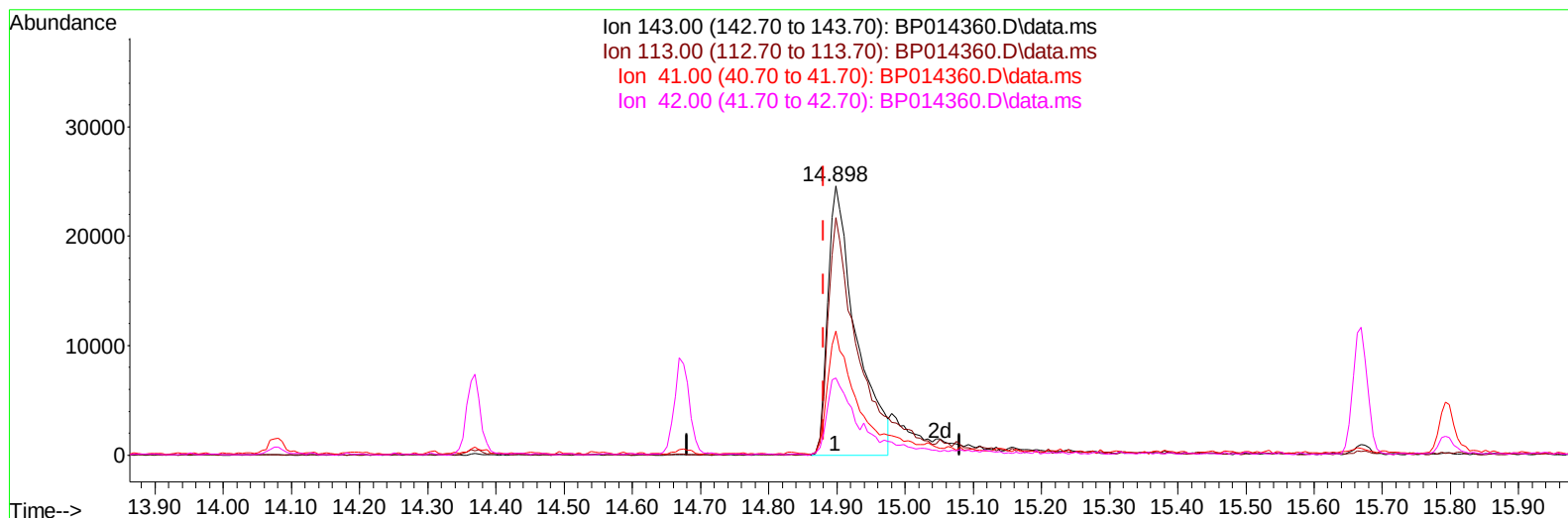
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032923\
 Data File : BP014360.D
 Acq On : 29 Mar 2023 13:08
 Operator : CG/JU
 Sample : 01982-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EYGC2

Manual IntegrationsAPPROVED

Quant Time: Mar 30 01:01:29 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 30 00:56:20 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/30/2023
 Supervised By :Jagrut Upadhyay 03/30/2023



TIC: BP014360.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.898min (+ 0.018) 14.61 ng/ul

response 69743

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	88.18
41.00	44.20	46.07
42.00	31.50	28.68

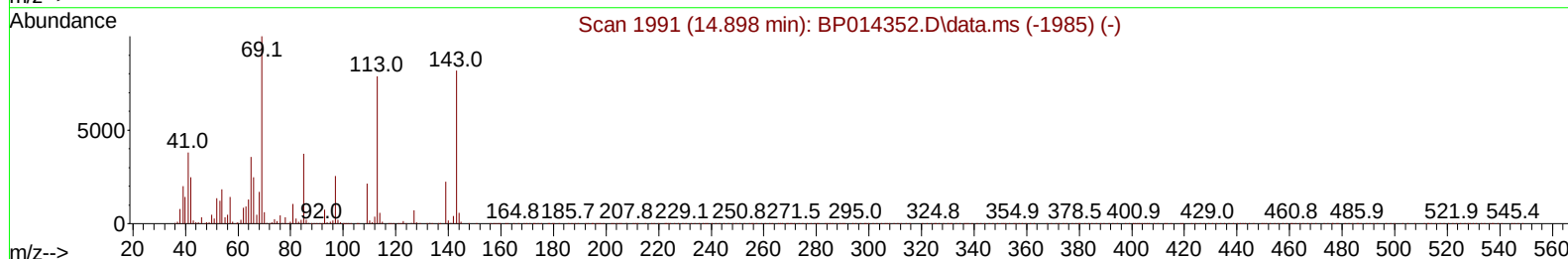
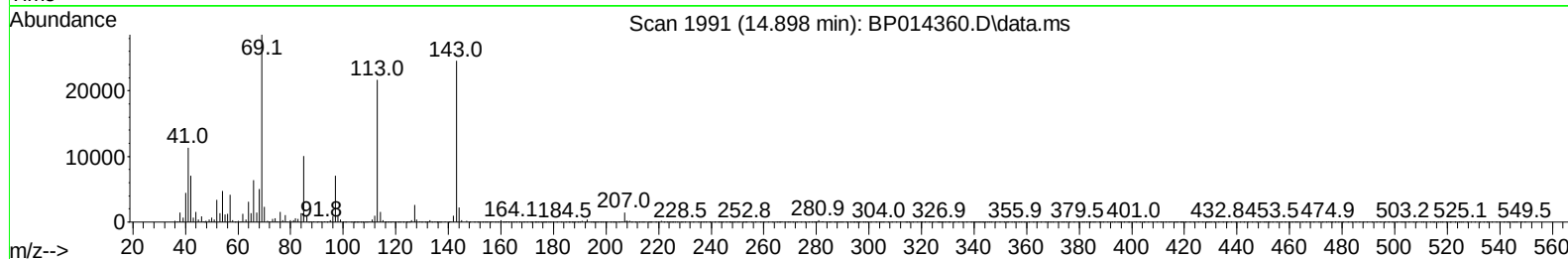
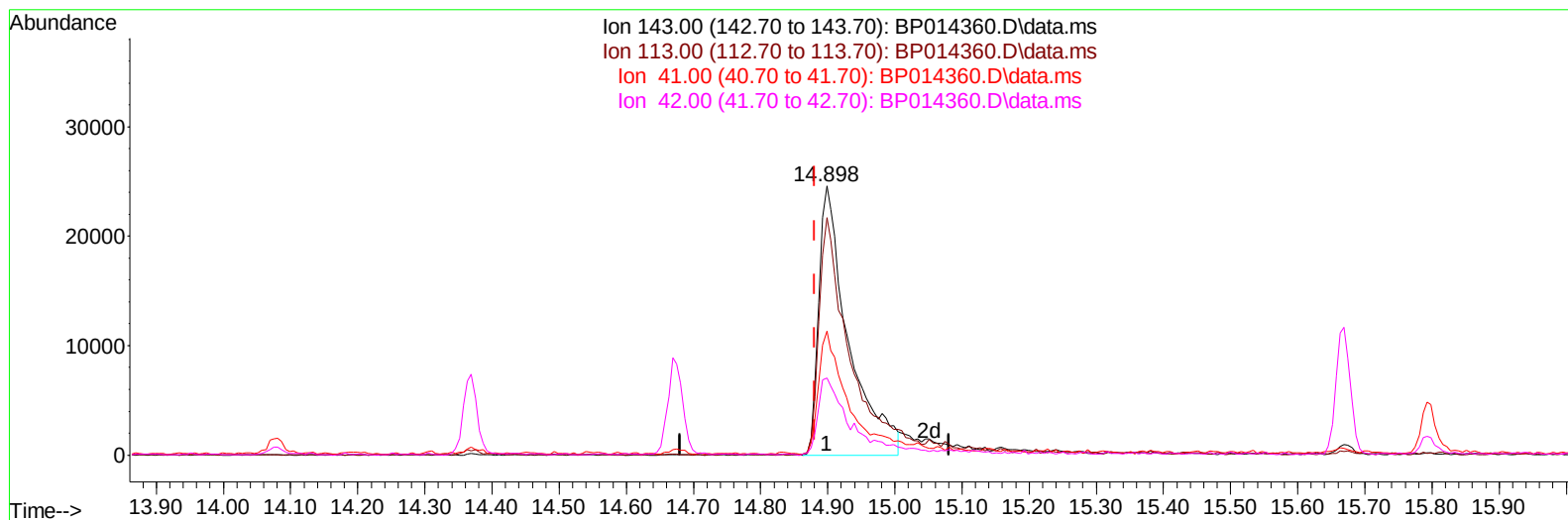
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(54) 4-Nitrophenol-d4 (S)

14.898min (+ 0.018) 15.69 ng/ul m

response 74927

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	88.18
41.00	44.20	46.07
42.00	31.50	28.68

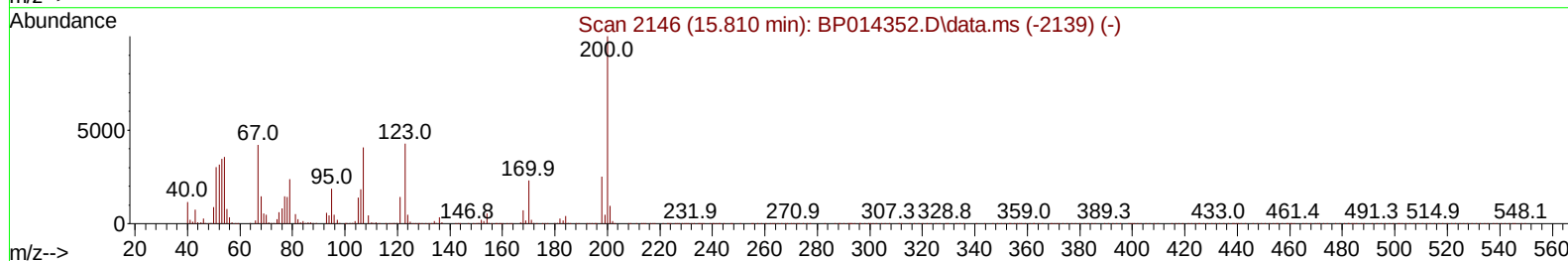
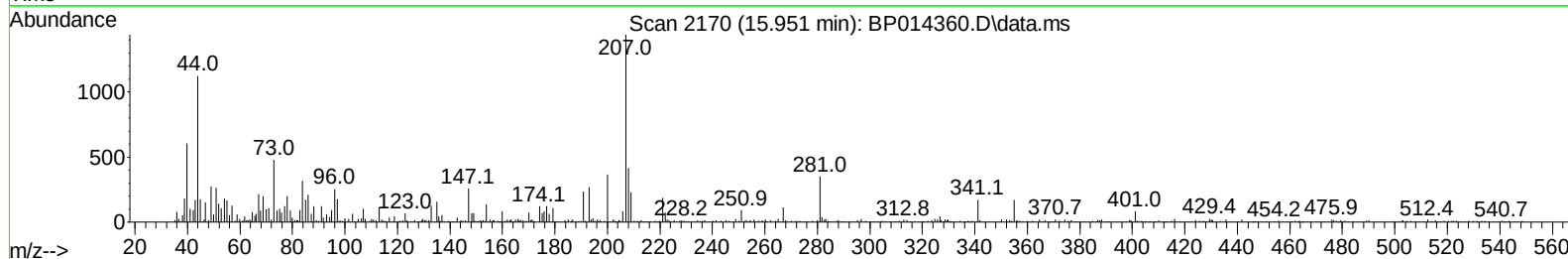
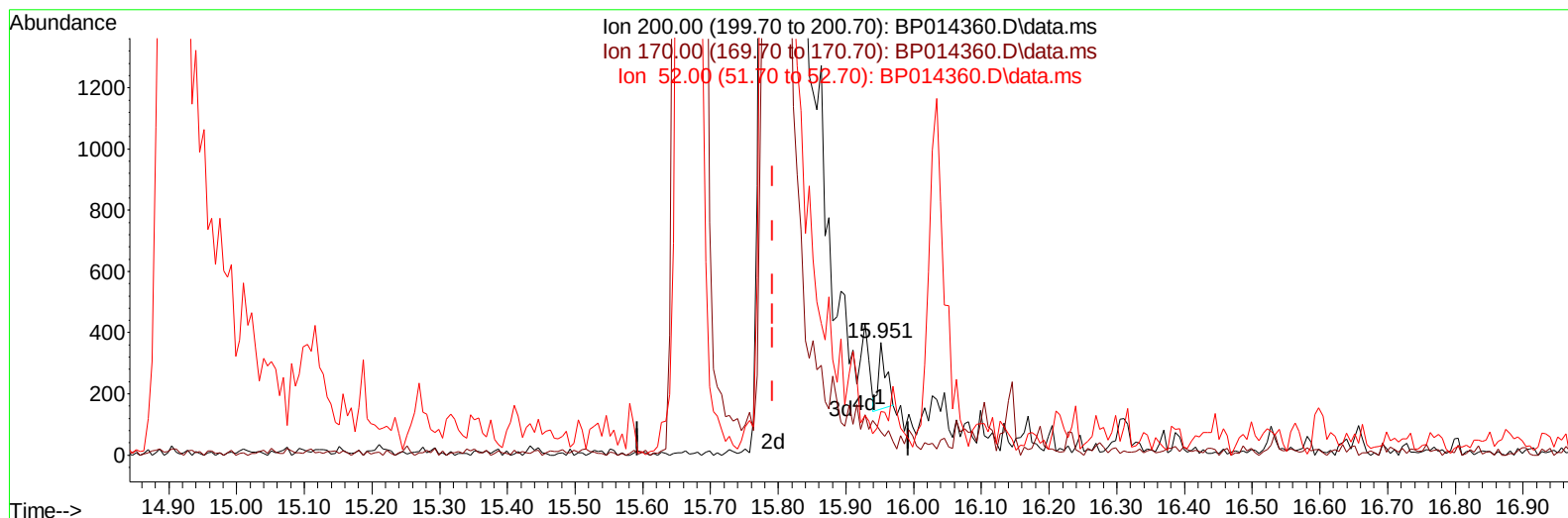
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TIC: BP014360.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.951min (+ 0.159) 0.03 ng/ul

response	173	
Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.10	20.44
52.00	45.50	38.96
0.00	0.00	0.00

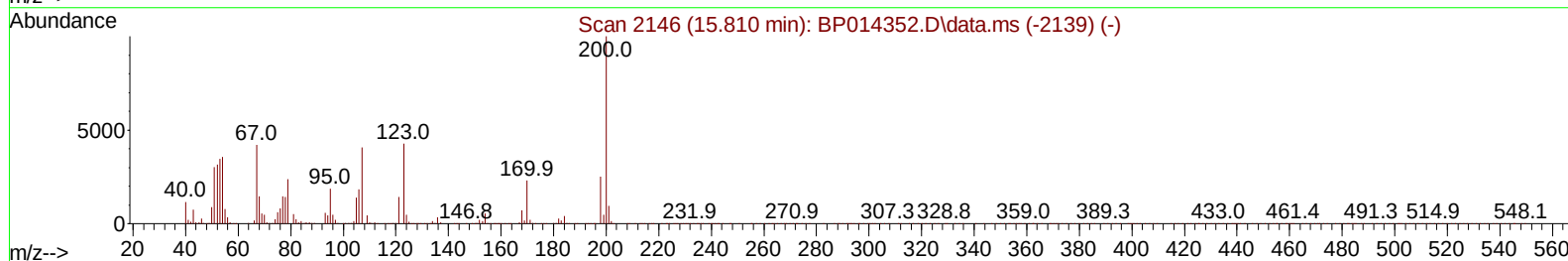
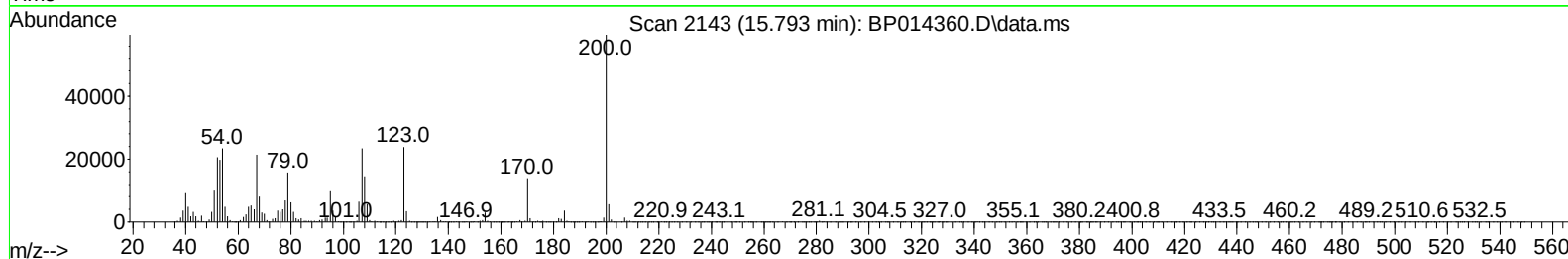
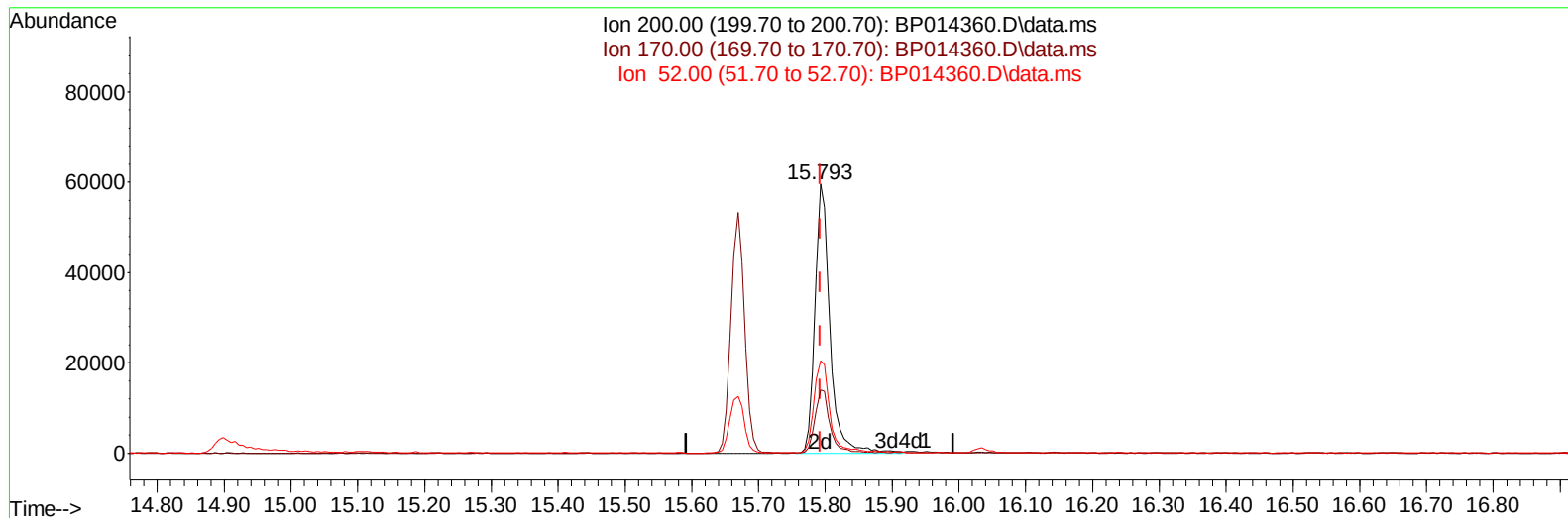
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Instrument :
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TIC: BP014360.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.793min (+ 0.000) 14.45 ng/ul m

response 92564

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.10	23.47
52.00	45.50	34.43#
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	8.022	152	132890	20.000	ng/ul	0.00
20) Naphthalene-d8	10.846	136	573238	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.675	164	389744	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.434	188	870599	20.000	ng/ul	0.00
79) Chrysene-d12	21.522	240	831472	20.000	ng/ul	0.00
88) Perylene-d12	24.051	264	772036	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	15010	4.380	ng/uL	0.00
4) Pyridine-d5	3.828	84	177540	20.217	ng/ul	0.00
7) Phenol-d5	7.181	99	258655	21.105	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.346	67	181805	23.121	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	198961	22.461	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	202735	20.508	ng/ul	0.00
21) Nitrobenzene-d5	9.199	128	100863	21.796	ng/ul	0.00
24) 2-Nitrophenol-d4	9.922	143	115124	21.726	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.469	165	194825	20.188	ng/ul	0.00
31) 4-Chloroaniline-d4	10.993	131	260178	20.468	ng/ul	0.00
46) Dimethylphthalate-d6	14.081	166	719504	22.768	ng/ul	0.00
49) Acenaphthylene-d8	14.369	160	788631	22.491	ng/ul	0.00
54) 4-Nitrophenol-d4	14.898	143	74927m	15.695	ng/ul	0.02
60) Fluorene-d10	15.669	176	625305	23.728	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.793	200	92564m	14.450	ng/ul	0.00
73) Anthracene-d10	17.534	188	966446	22.938	ng/ul	0.00
81) Pyrene-d10	19.763	212	1162307	20.811	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.880	264	979780	23.887	ng/ul	0.00

Target Compounds Qvalue

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

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