

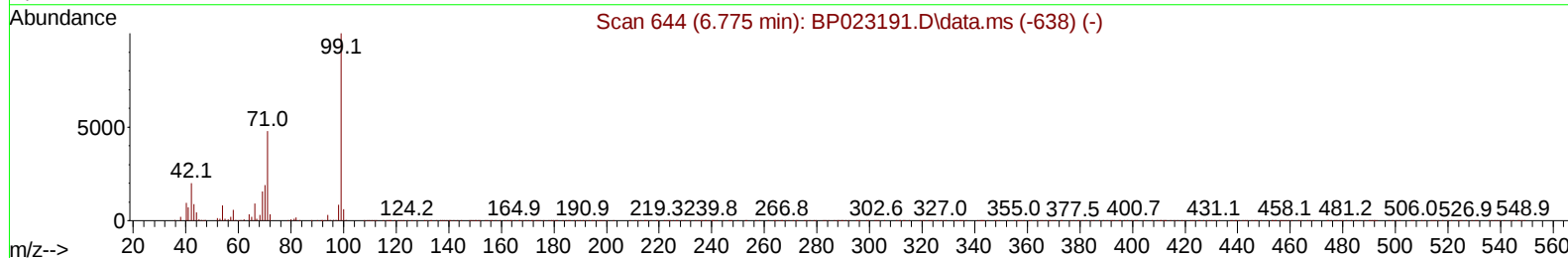
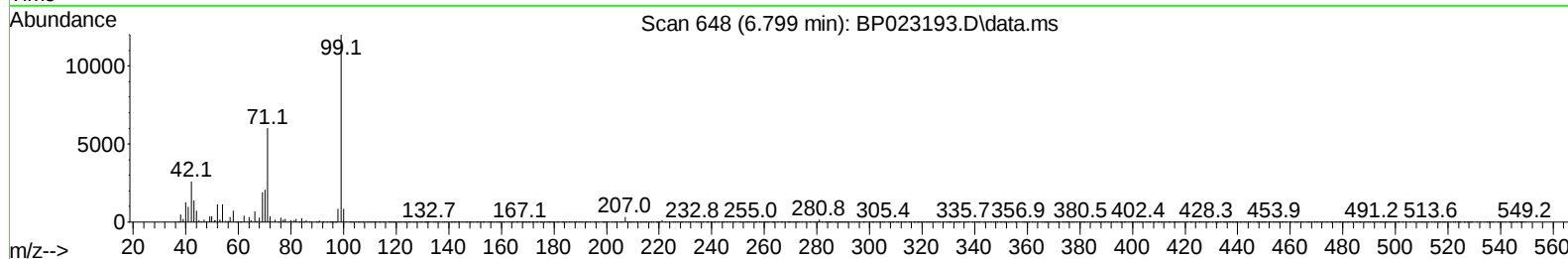
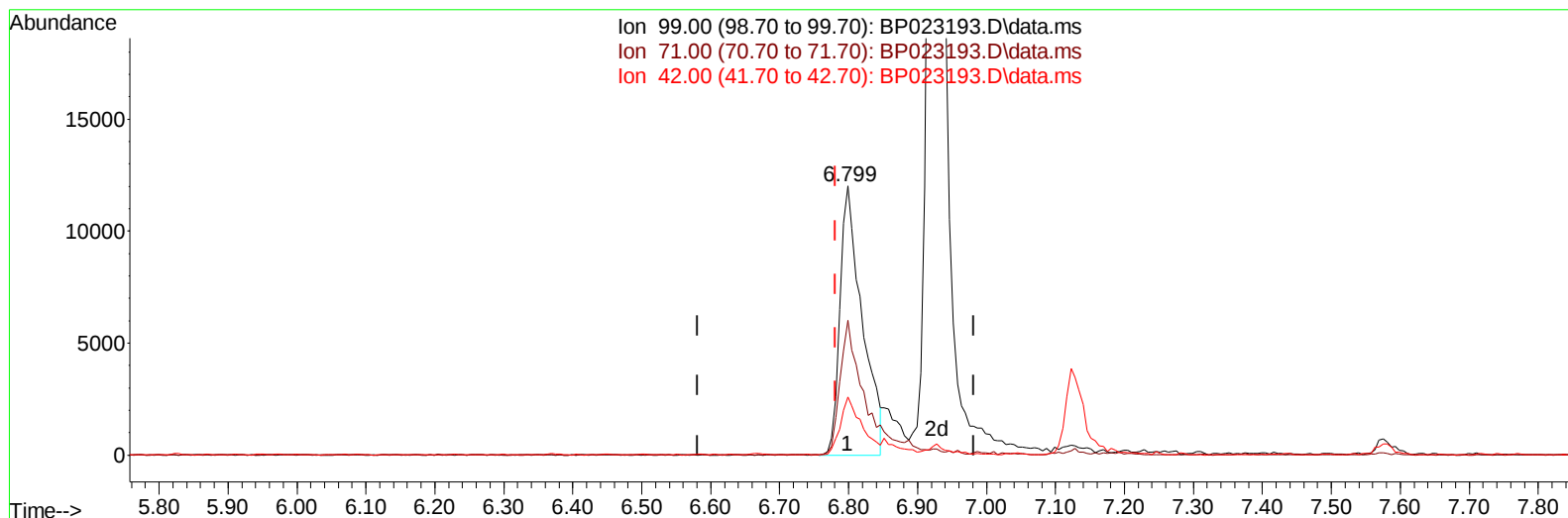
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111824\
Data File : BP023193.D
Acq On : 18 Nov 2024 11:49
Operator : RC/JU
Sample : P4829-14
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E29S0

Manual IntegrationsAPPROVED

Quant Time: Nov 18 12:25:46 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 13 04:49:01 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/19/2024
Supervised By :mohammad ahmed 11/19/2024



TIC: BP023193.D\data.ms

(7) Phenol-d5 (S)

6.799min (+ 0.018) 7.27 ng/u1

response 26744

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	44.70	50.15
42.00	21.10	21.56
0.00	0.00	0.00

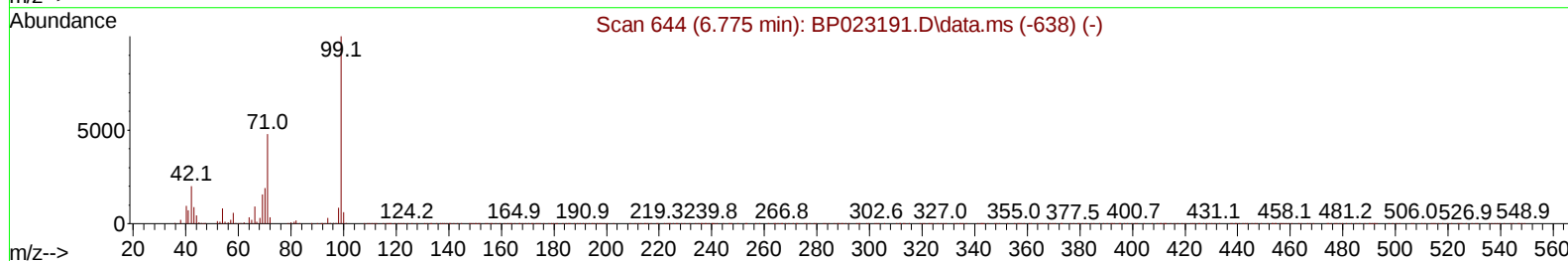
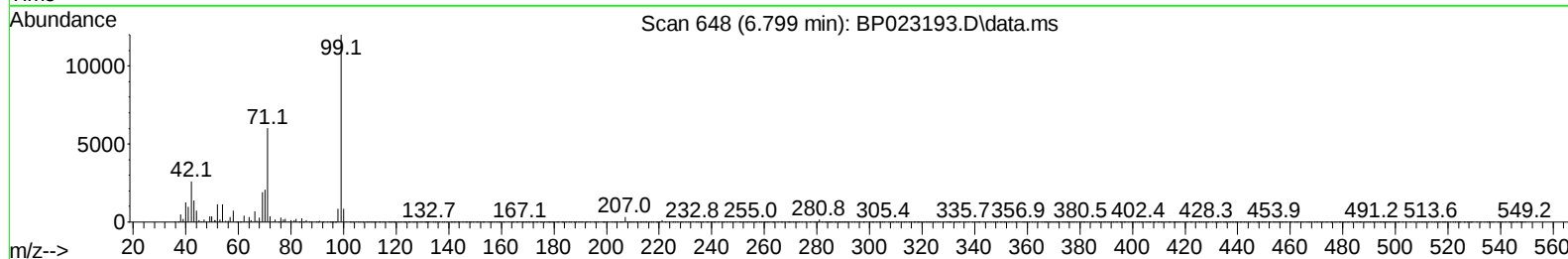
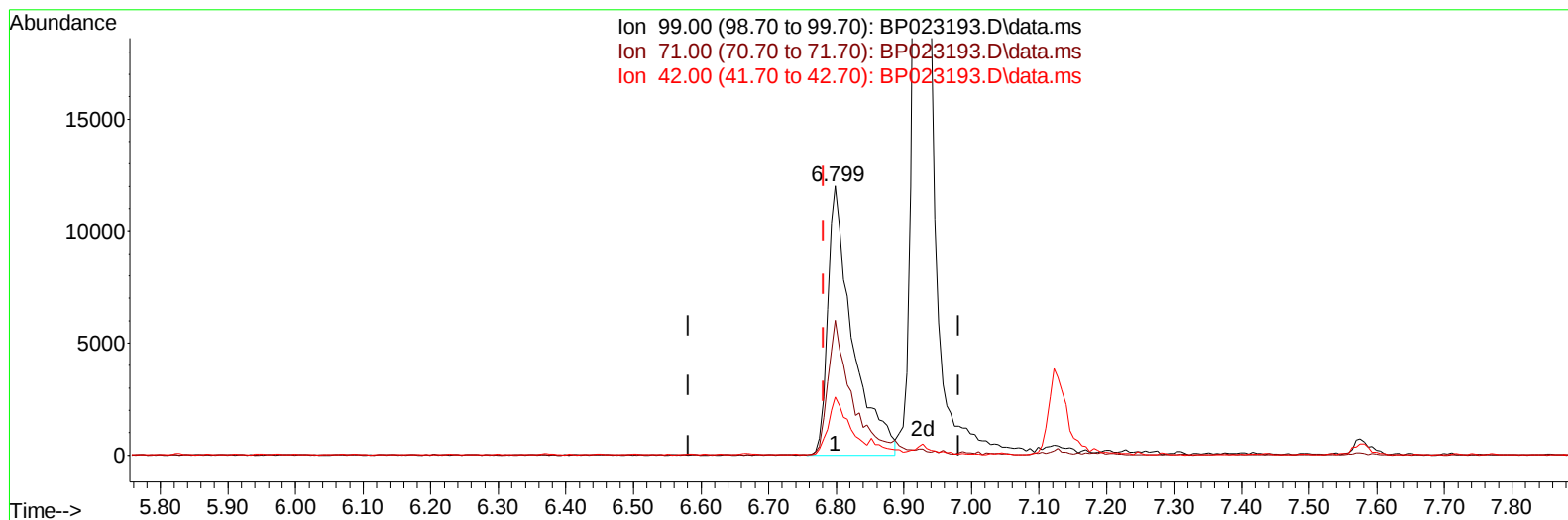
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(7) Phenol-d5 (S)

6.799min (+ 0.018) 8.24 ng/ul m

response 30344

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	44.70	50.15
42.00	21.10	21.56
0.00	0.00	0.00

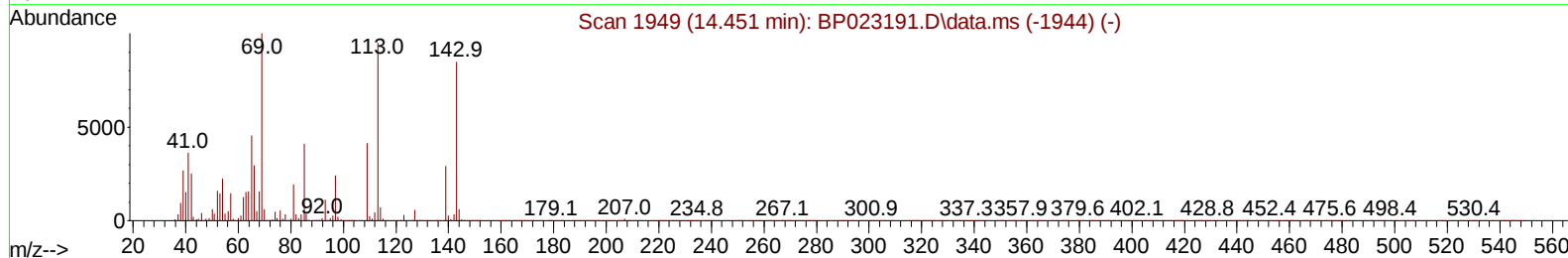
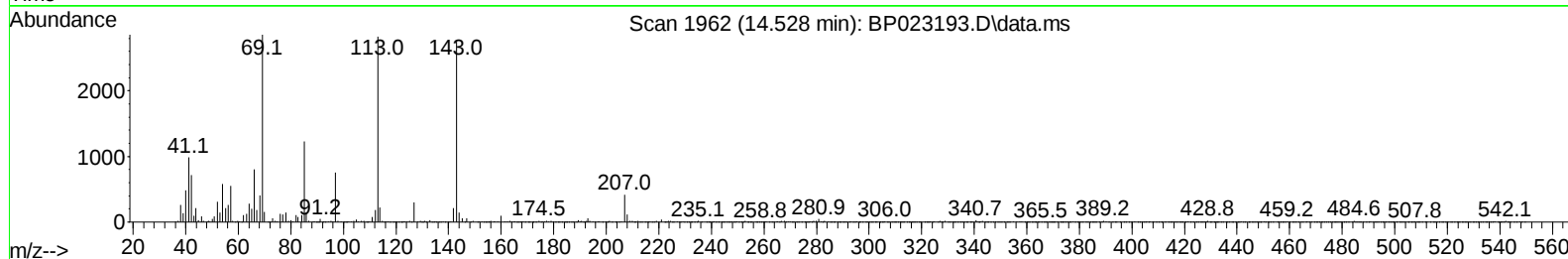
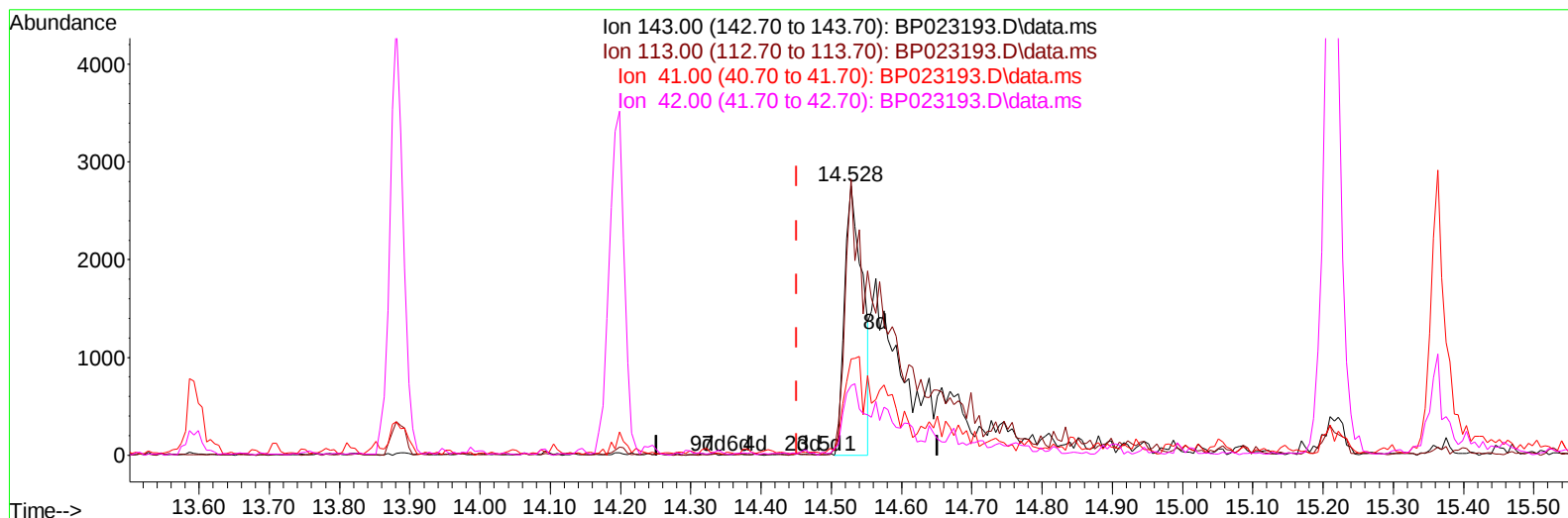
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TIC: BP023193.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.076) 2.91 ng/u1

response 4896

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	102.72
41.00	39.20	35.74
42.00	25.90	25.90

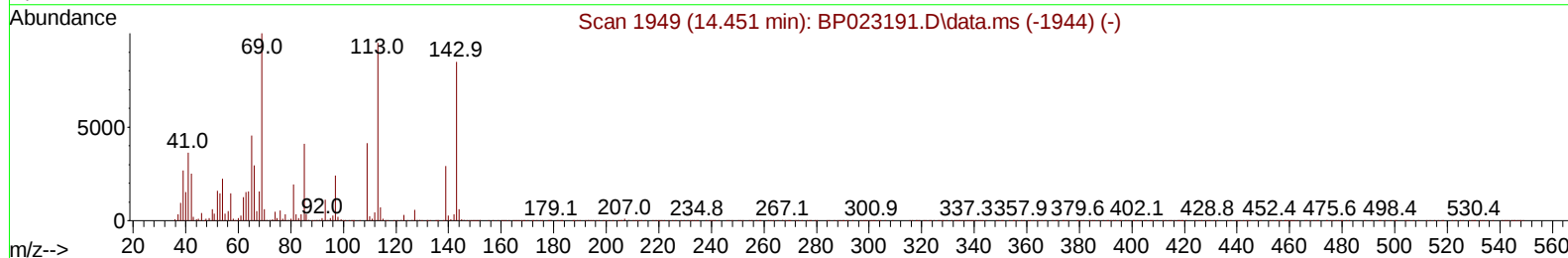
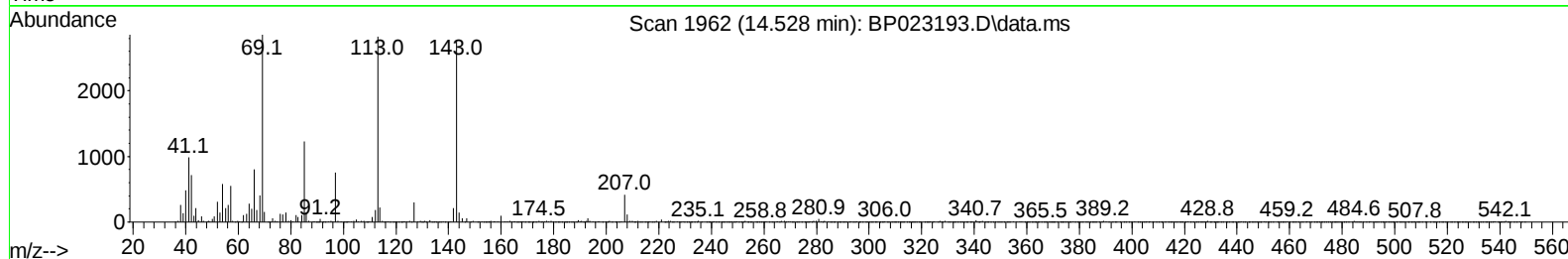
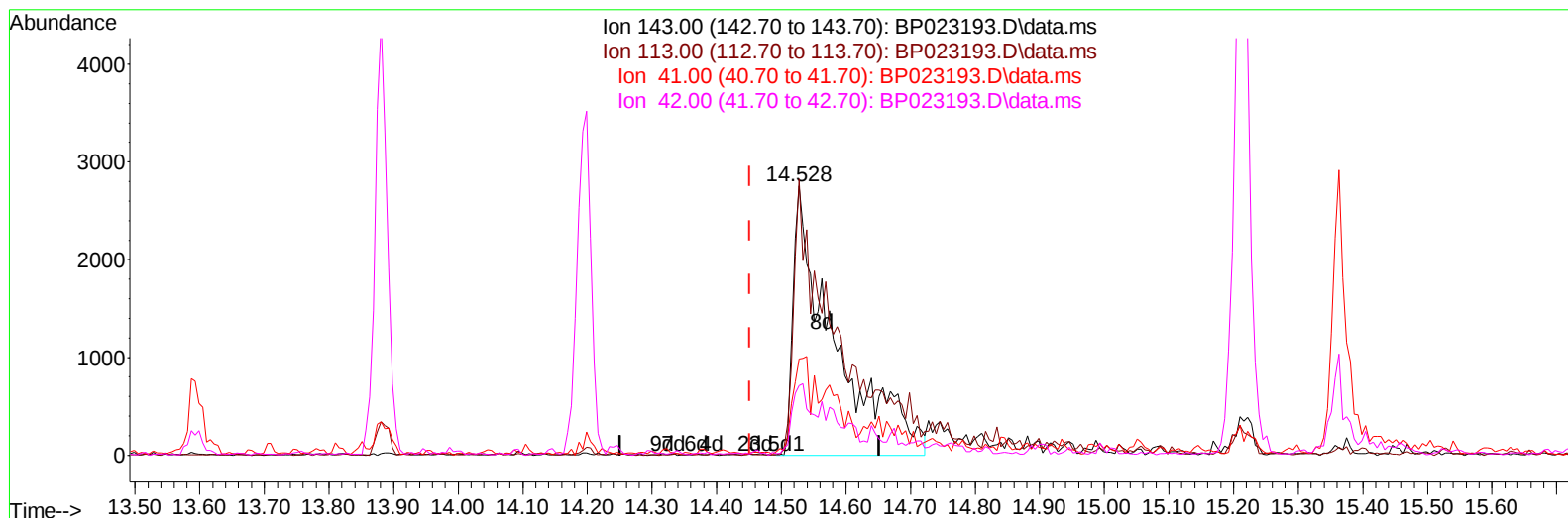
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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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TIC: BP023193.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.076) 7.27 ng/ul m

response 12242

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	102.72
41.00	39.20	35.74
42.00	25.90	25.90

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 ALS Vial : 4 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/19/2024
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Quant Time: Nov 18 13:24:27 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	49786	20.000	ng/ul	0.00
20) Naphthalene-d8	10.322	136	192846	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.198	164	155405	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.004	188	386335	20.000	ng/ul	-0.01
79) Chrysene-d12	21.445	240	440456	20.000	ng/ul	0.00
88) Perylene-d12	24.668	264	511226	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	5891	5.688	ng/uL	0.00
4) Pyridine-d5	3.564	84	19078	6.200	ng/ul	0.02
7) Phenol-d5	6.799	99	30344m	8.244	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	6.922	67	71149	32.900	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	79869	29.628	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	59830	19.752	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	44697	33.293	ng/ul	0.00
24) 2-Nitrophenol-d4	9.428	143	52187	32.606	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	104594	31.205	ng/ul	0.01
31) 4-Chloroaniline-d4	10.481	131	101863	25.474	ng/ul	0.00
46) Dimethylphthalate-d6	13.592	166	399928	35.706	ng/ul	-0.02
49) Acenaphthylene-d8	13.881	160	402278	34.421	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.528	143	12242m	7.273	ng/ul	0.08
60) Fluorene-d10	15.216	176	363298	36.996	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.363	200	66295	28.645	ng/ul	0.00
73) Anthracene-d10	17.104	188	695181	42.498	ng/ul	-0.01
81) Pyrene-d10	19.492	212	897750	44.251	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.439	264	1053811	43.275	ng/ul	0.00

Target Compounds Qvalue

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

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