

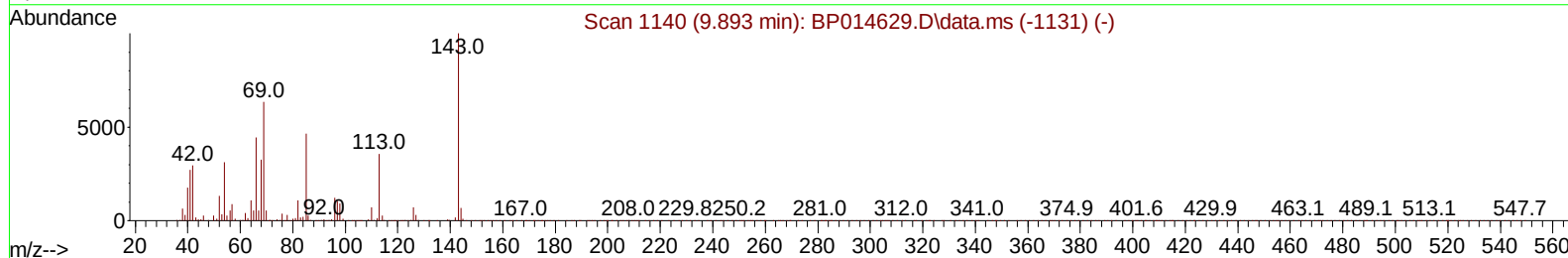
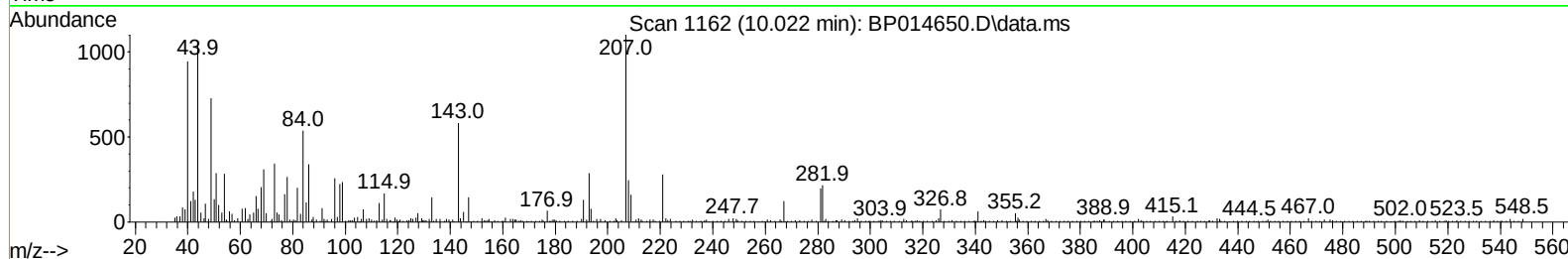
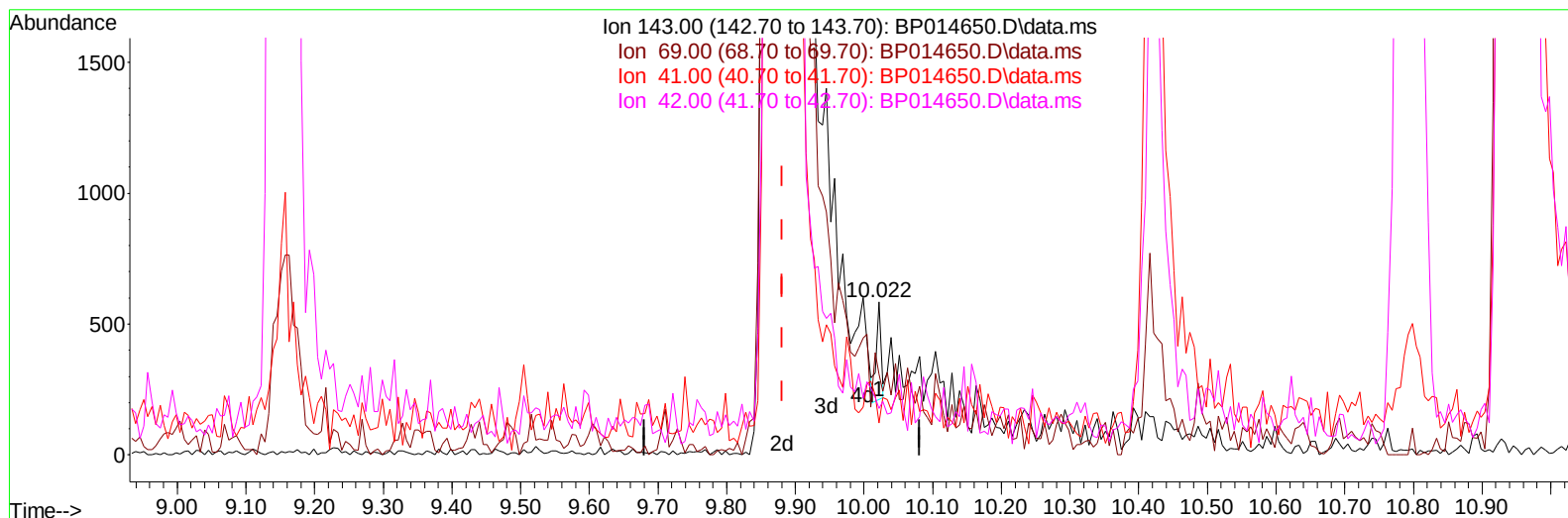
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041023\
 Data File : BP014650.D
 Acq On : 11 Apr 2023 11:44
 Operator : CG/JU
 Sample : PB152027BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK027

Manual IntegrationsAPPROVED

Quant Time: Apr 11 23:00:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 11 01:21:43 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/12/2023
 Supervised By :Jagrut Upadhyay 04/12/2023



TIC: BP014650.D\data.ms

(24) 2-Nitrophenol-d4 (S)

10.022min (+ 0.141) 0.03 ng/ul

response 192

Ion	Exp%	Act%
143.00	100.00	100.00
69.00	62.40	52.91
41.00	24.10	20.89
42.00	25.90	30.48

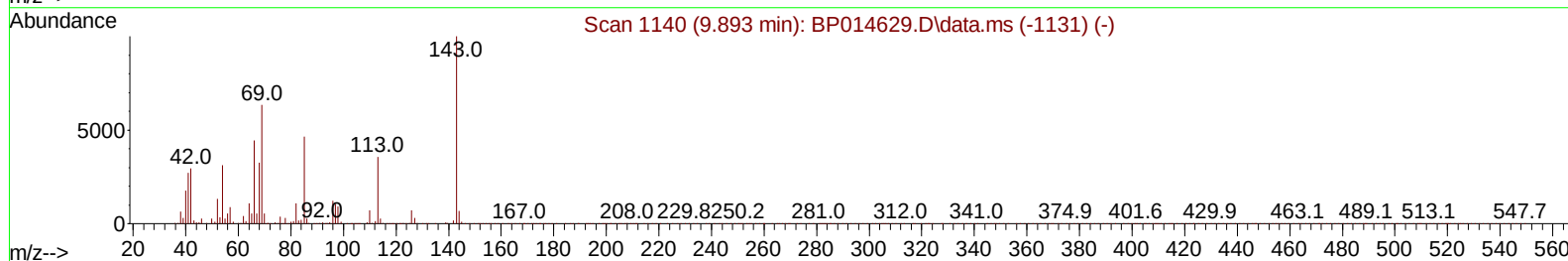
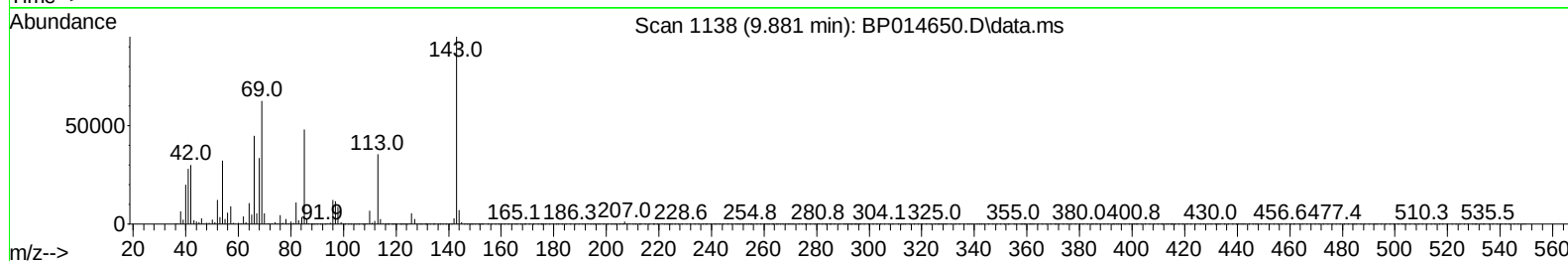
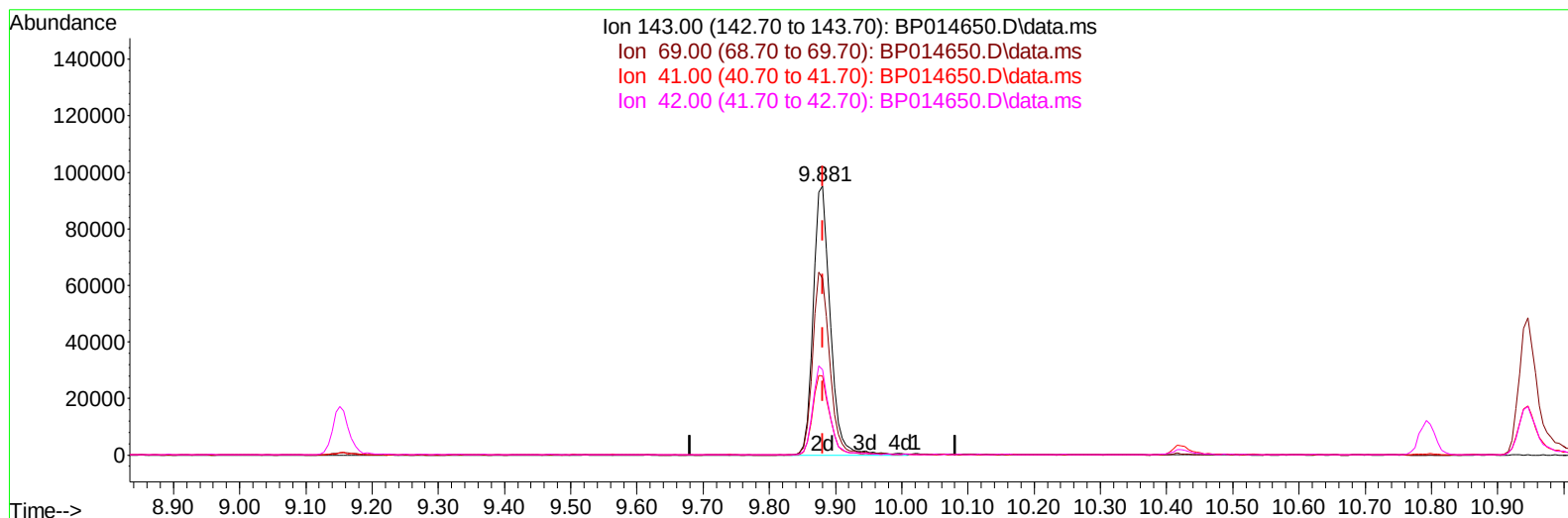
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(24) 2-Nitrophenol-d4 (S)

9.881min (-0.000) 29.46 ng/ul m

response 170775

Ion	Exp%	Act%
143.00	100.00	100.00
69.00	62.40	65.71
41.00	24.10	29.43#
42.00	25.90	31.65#

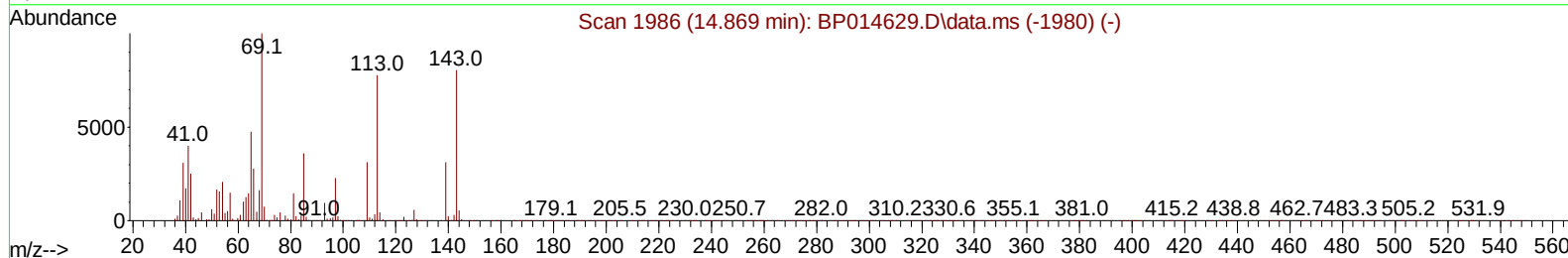
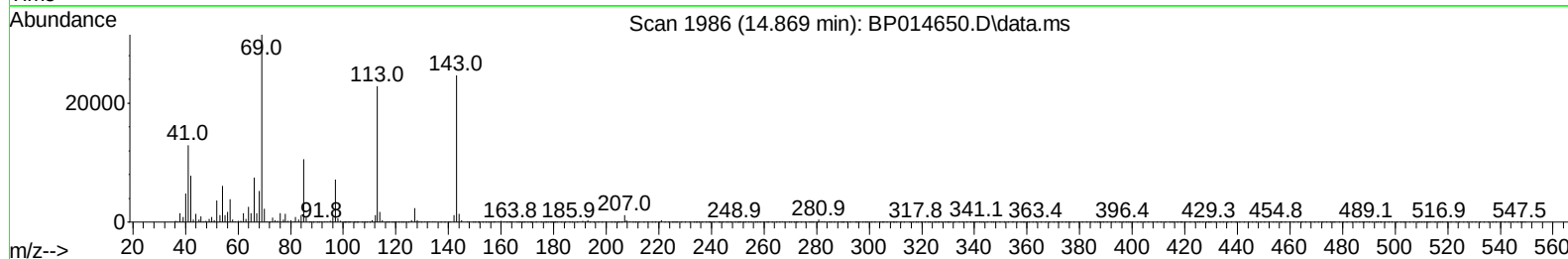
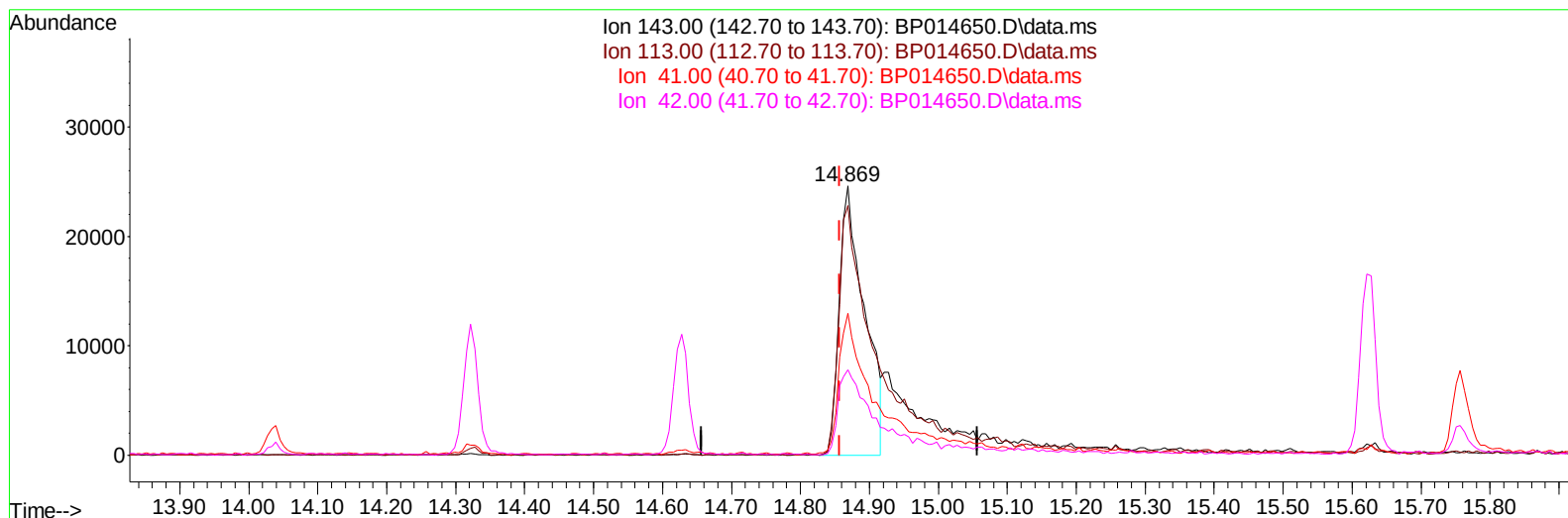
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TIC: BP014650.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.869min (+ 0.011) 13.22 ng/ul

response 62144

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	92.78
41.00	44.20	52.61
42.00	31.50	31.83

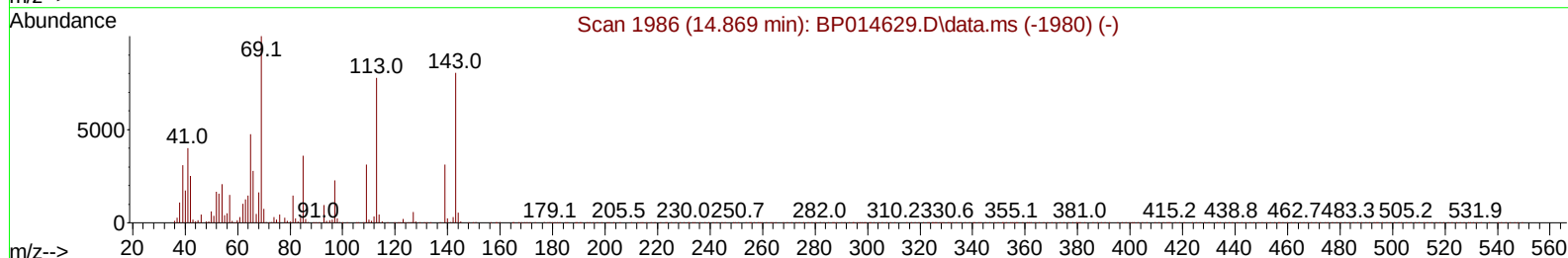
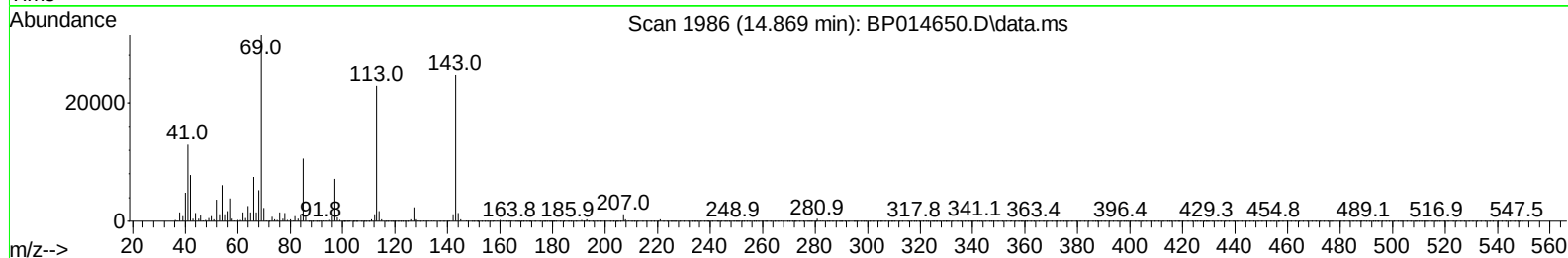
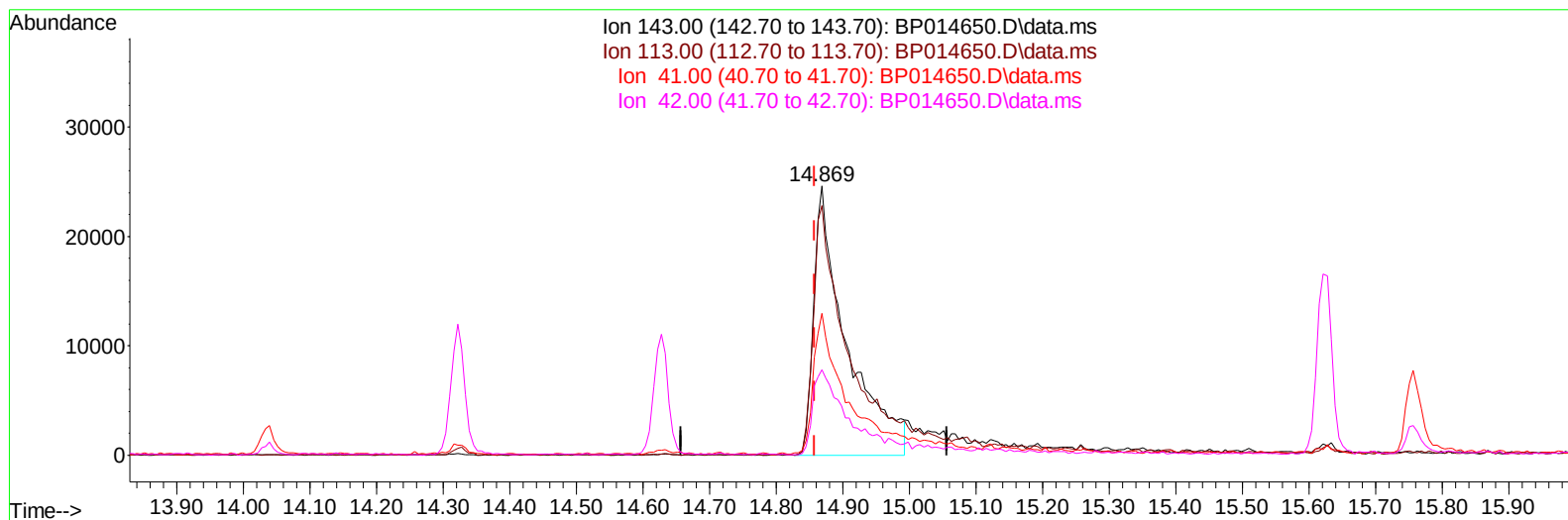
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Sample : PB152027BL
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
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TIC: BP014650.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.869min (+ 0.011) 17.86 ng/ul m

response 83902

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	92.78
41.00	44.20	52.61
42.00	31.50	31.83

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	142278	20.000	ng/ul	0.00
20) Naphthalene-d8	10.792	136	627057	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.627	164	383620	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	745906	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	569520	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	688955	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	21923	5.975	ng/uL	0.00
4) Pyridine-d5	3.793	84	282786	30.077	ng/ul	0.00
7) Phenol-d5	7.140	99	383052	29.193	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.304	67	277807	32.998	ng/ul	0.00
11) 2-Chlorophenol-d4	7.504	132	293687	30.967	ng/ul	0.00
15) 4-Methylphenol-d8	8.692	113	314077	29.674	ng/ul	0.00
21) Nitrobenzene-d5	9.151	128	146979	29.035	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	170775m	29.462	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.422	165	273688	25.926	ng/ul	0.00
31) 4-Chloroaniline-d4	10.945	131	431241	31.013	ng/ul	0.00
46) Dimethylphthalate-d6	14.039	166	948689	30.500	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	1052217	30.487	ng/ul	0.00
54) 4-Nitrophenol-d4	14.869	143	83902m	17.855	ng/ul	0.01
60) Fluorene-d10	15.627	176	766562	29.553	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	123934	22.581	ng/ul	0.00
73) Anthracene-d10	17.492	188	1148645	31.820	ng/ul	0.00
81) Pyrene-d10	19.721	212	1189856	31.104	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1217950	33.274	ng/ul	0.00

Target Compounds Qvalue

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

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