

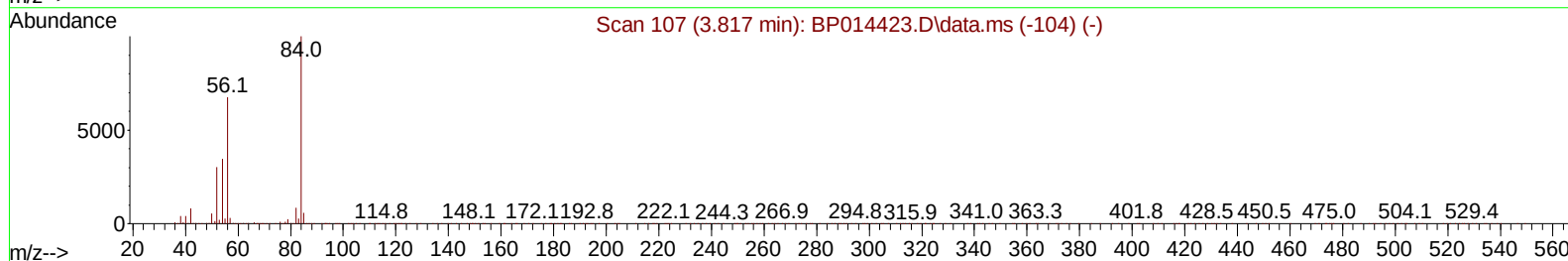
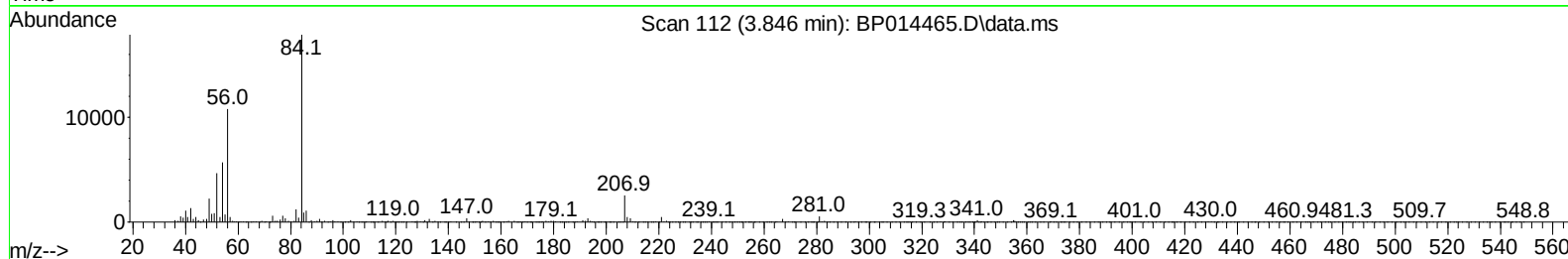
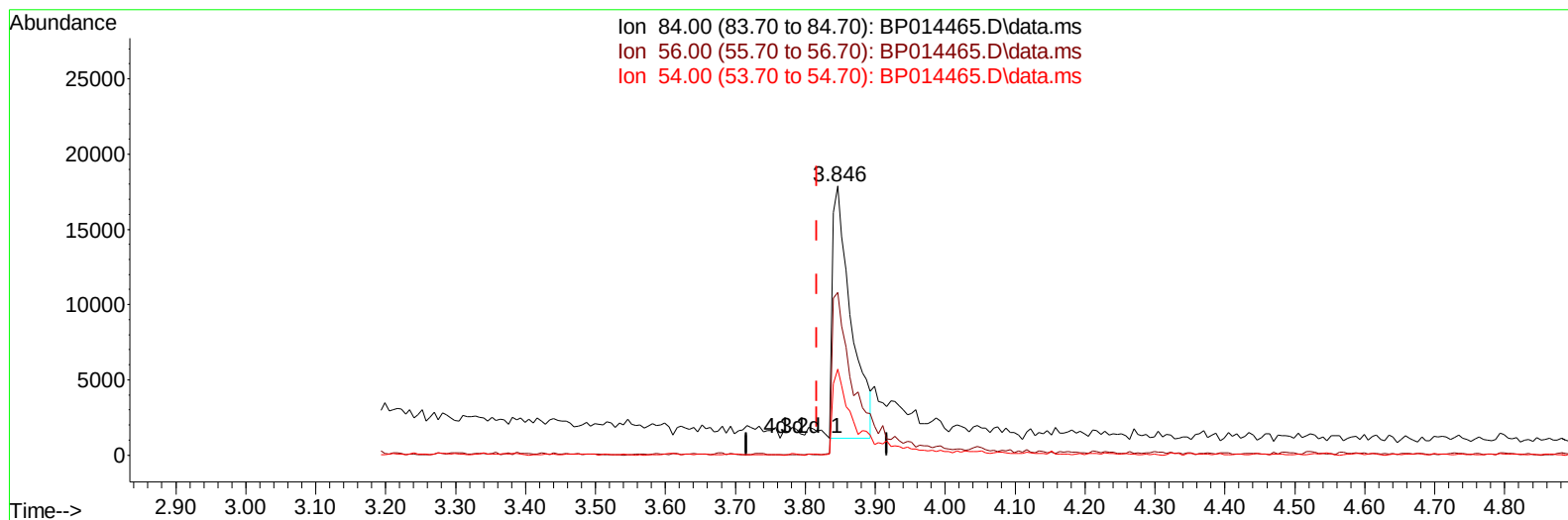
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014465.D
Acq On : 01 Apr 2023 03:48
Operator : CG/JU
Sample : 02092-14
Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB981

Manual IntegrationsAPPROVED

Quant Time: Apr 01 04:44:07 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 31 04:41:29 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/03/2023
Supervised By :Jagrut Upadhyay 04/03/2023



TIC: BP014465.D\data.ms

(4) Pyridine-d5 (S)

3.846min (+ 0.029) 2.82 ng/u1

response 30874

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	60.51
54.00	32.60	31.89
0.00	0.00	0.00

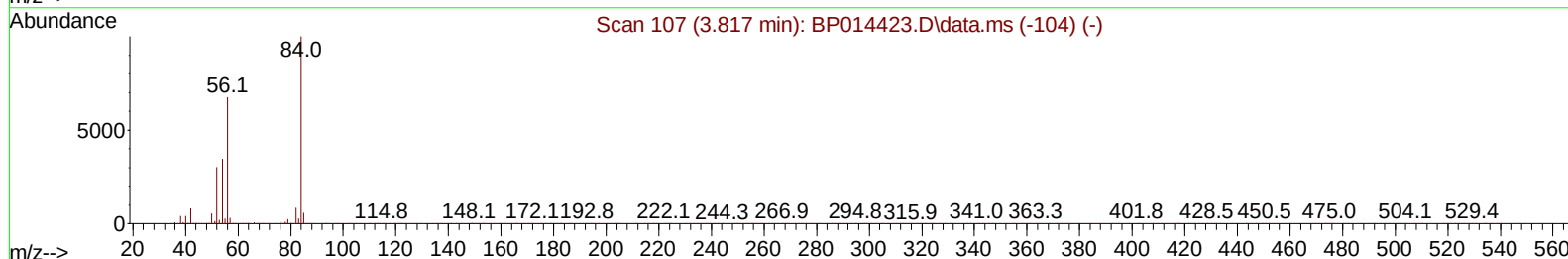
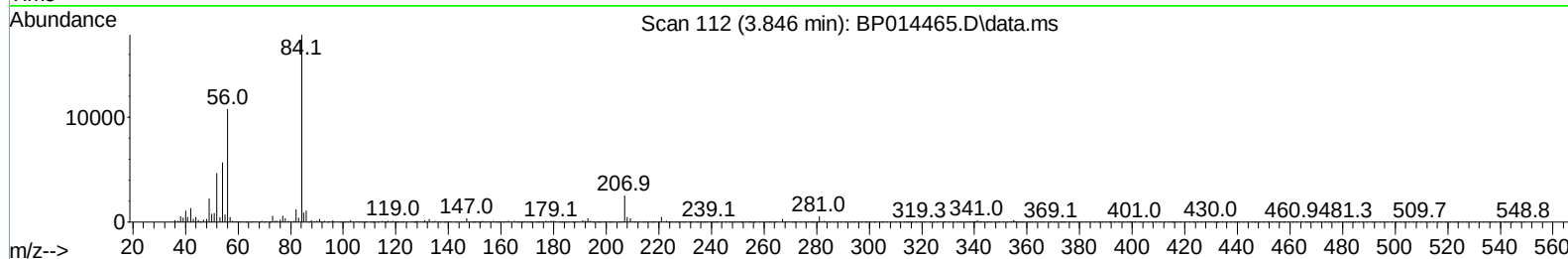
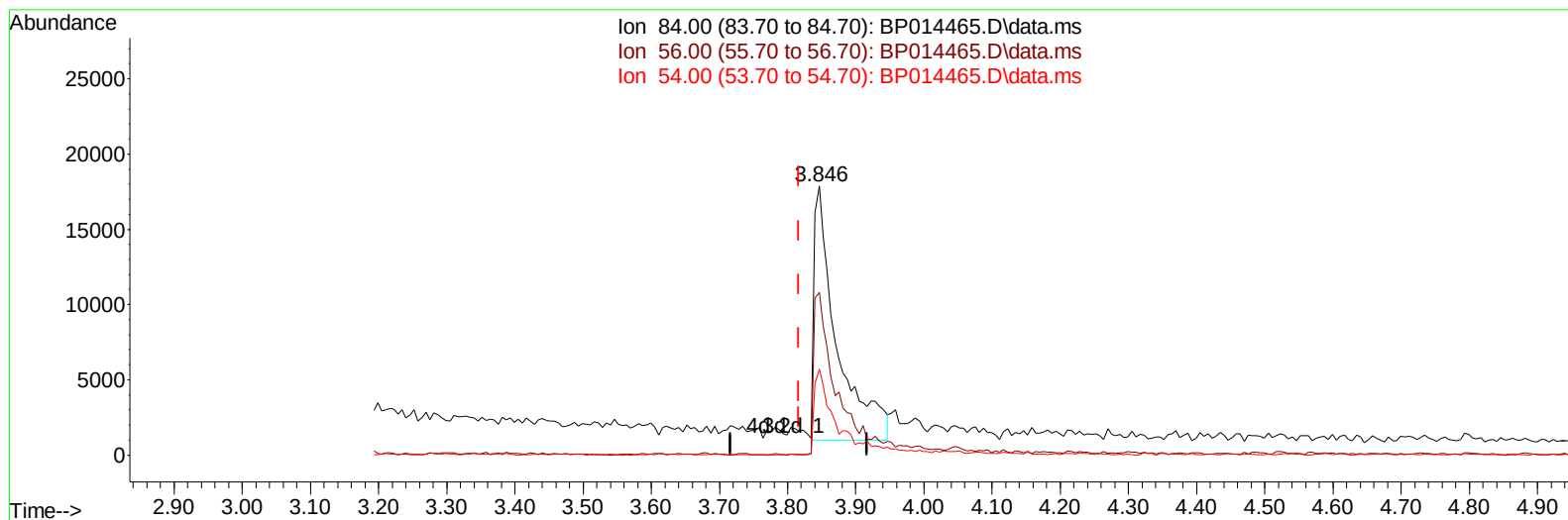
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(4) Pyridine-d5 (S)

3.846min (+ 0.029) 3.58 ng/ul m

response 39211

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	60.51
54.00	32.60	31.89
0.00	0.00	0.00

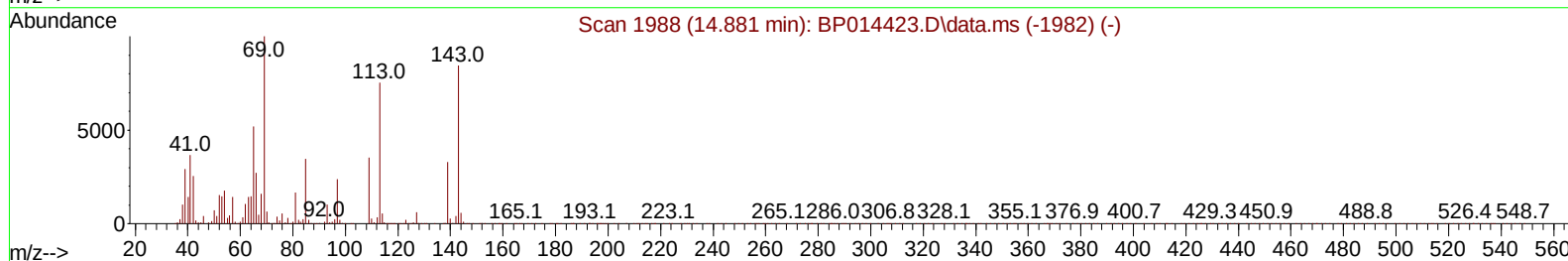
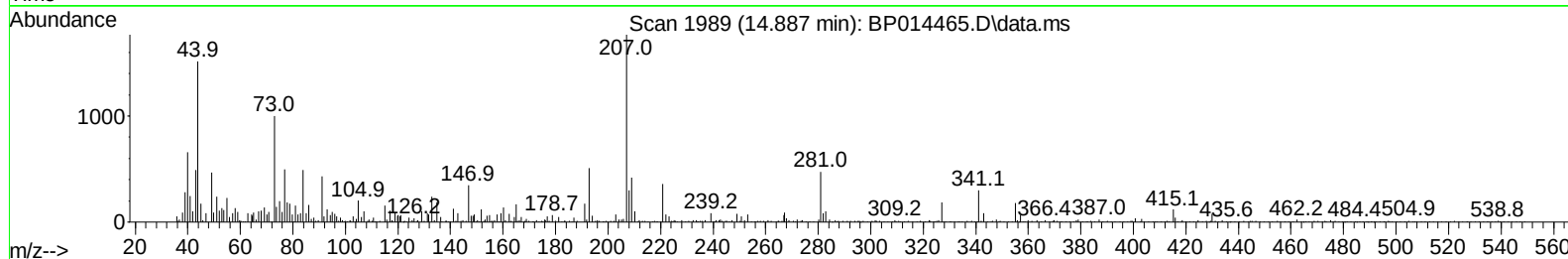
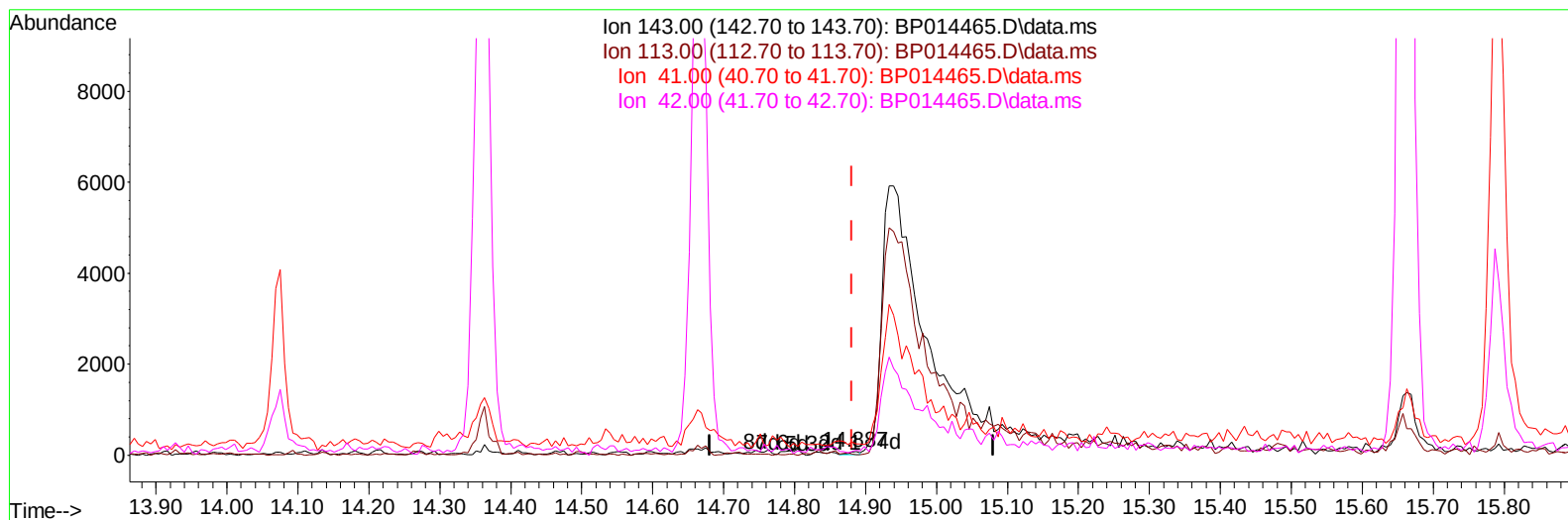
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TIC: BP014465.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.887min (+ 0.006) 0.01 ng/ul

response 60

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	17.07#
41.00	44.20	296.34#
42.00	31.50	123.17#

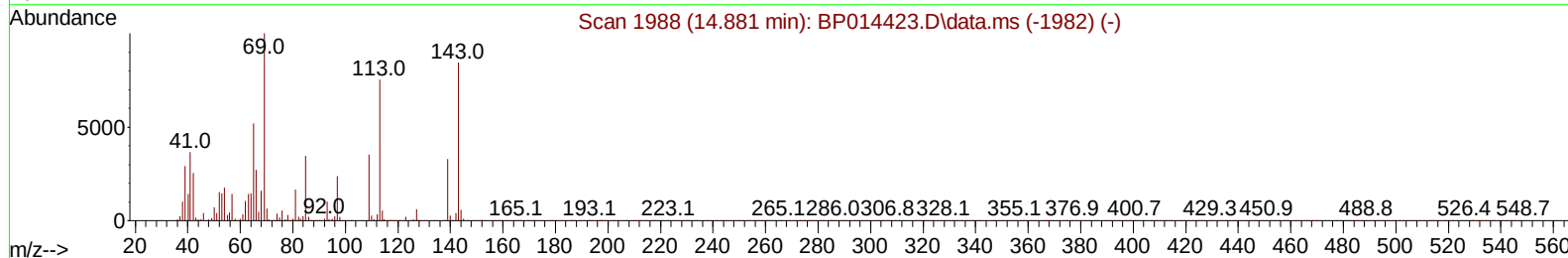
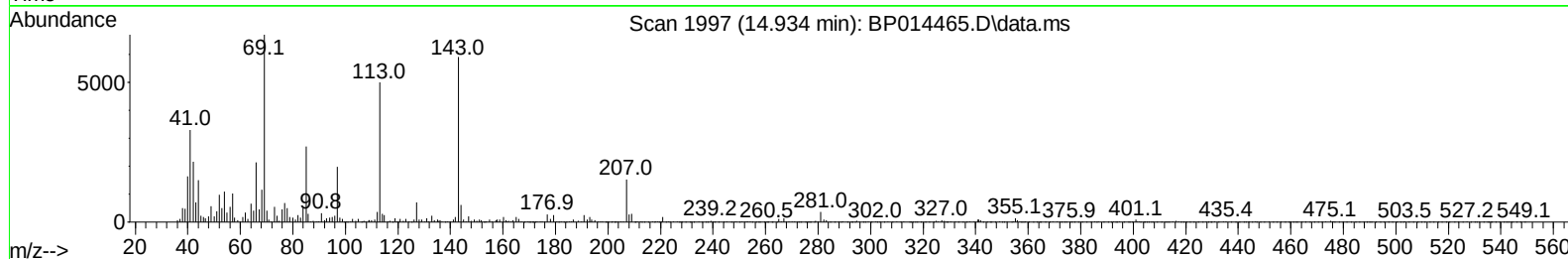
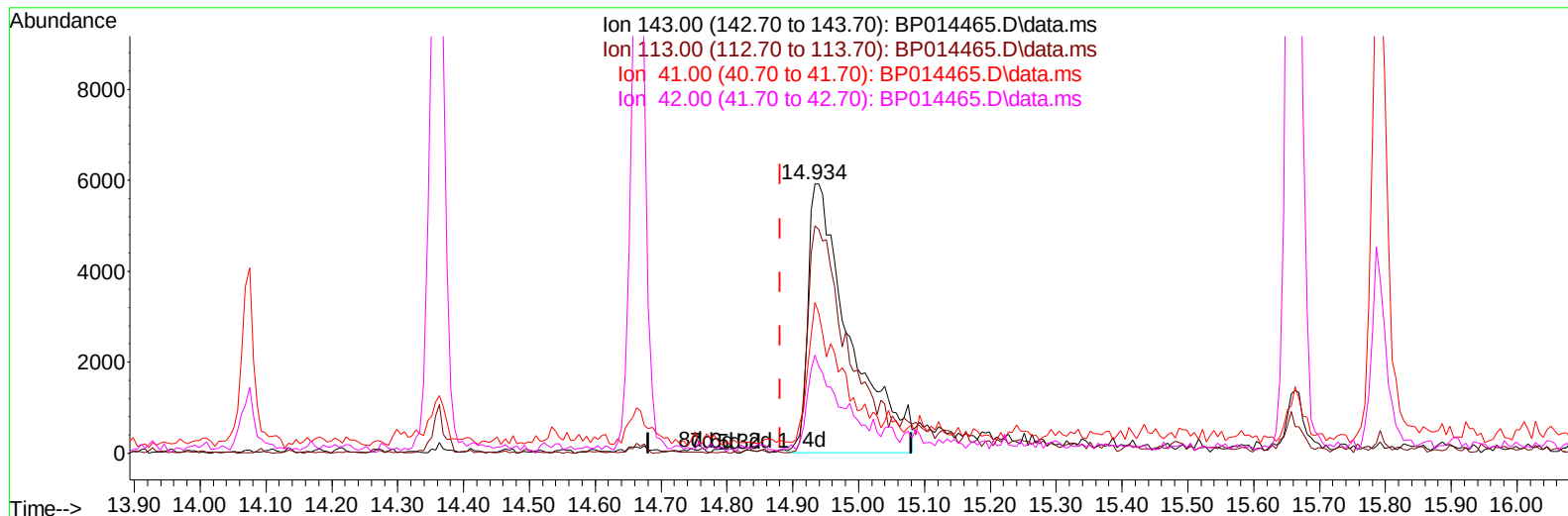
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Misc :
ALS Vial : 75 Sample Multiplier: 1

Instrument :
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TIC: BP014465.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.934min (+ 0.053) 4.01 ng/ul m

response 25993

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	84.62
41.00	44.20	55.98#
42.00	31.50	36.52

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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	8.011	152	165514	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	756897	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	529234	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	1157051	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	834195	20.000	ng/ul	0.00
88) Perylene-d12	24.039	264	807705	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	17240	4.039	ng/uL	0.00
4) Pyridine-d5	3.846	84	39211m	3.585	ng/ul	0.03
7) Phenol-d5	7.187	99	89974	5.894	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.334	67	311148	31.770	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	265504	24.065	ng/ul	0.00
15) 4-Methylphenol-d8	8.740	113	185853	15.094	ng/ul	0.01
21) Nitrobenzene-d5	9.187	128	186269	30.484	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	212869	30.424	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	352563	27.669	ng/ul	0.00
31) 4-Chloroaniline-d4	10.993	131	63398	3.777	ng/ul	0.01
46) Dimethylphthalate-d6	14.075	166	1405873	32.762	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	1503936	31.586	ng/ul	0.00
54) 4-Nitrophenol-d4	14.934	143	25993m	4.010	ng/ul	0.05
60) Fluorene-d10	15.663	176	1178997	32.947	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	205153	24.097	ng/ul	0.00
73) Anthracene-d10	17.528	188	1995735	35.641	ng/ul	0.00
81) Pyrene-d10	19.757	212	2158777	38.527	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.874	264	1557011	36.283	ng/ul	0.00

Target Compounds Qvalue

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

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