

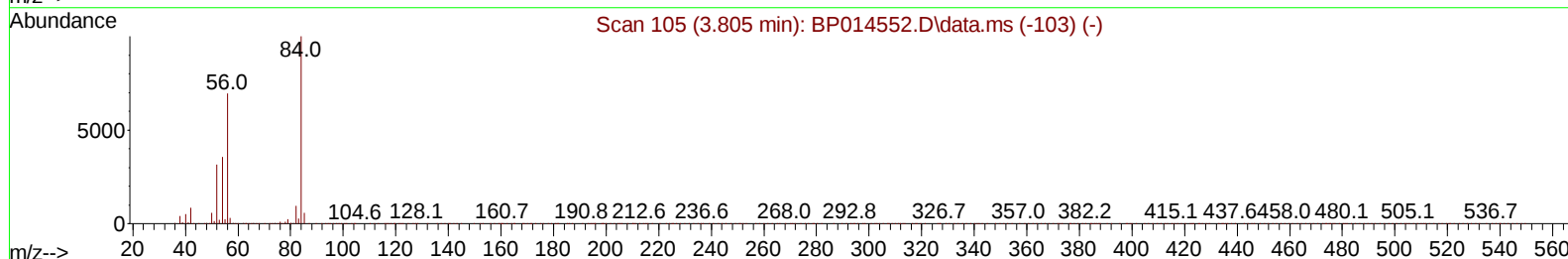
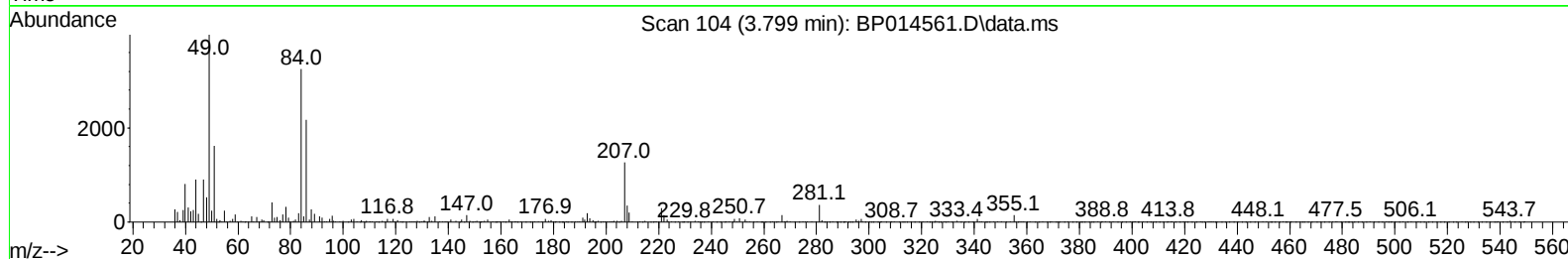
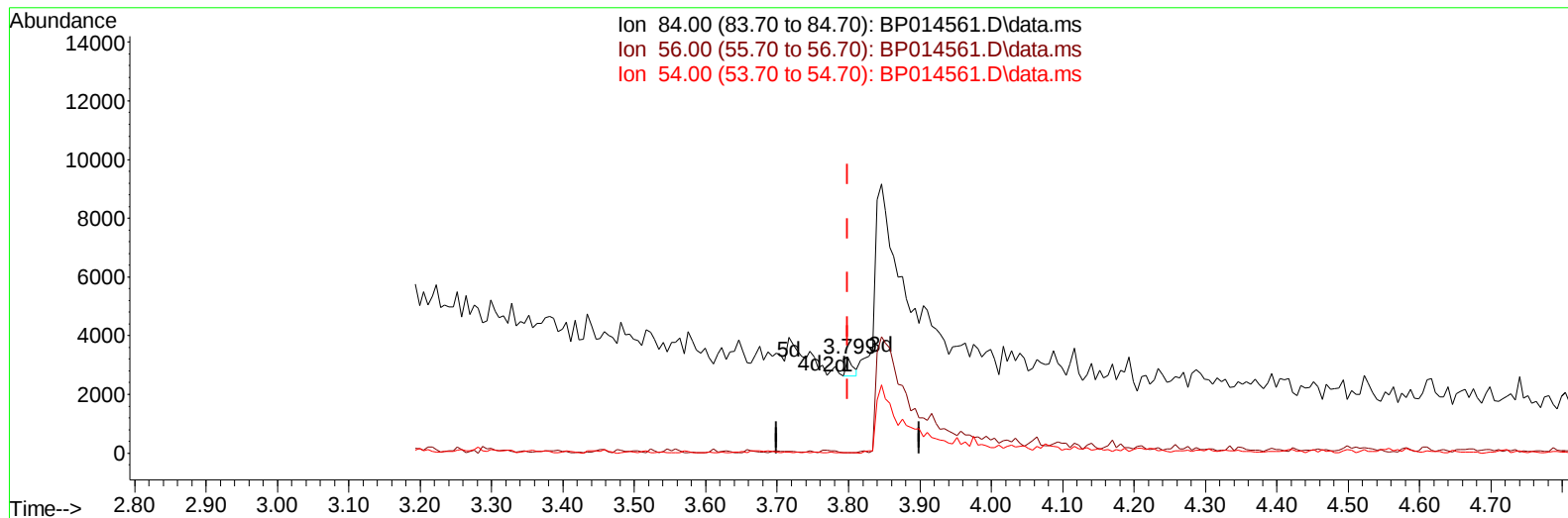
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
Data File : BP014561.D
Acq On : 05 Apr 2023 17:10
Operator : CG/JU
Sample : 02122-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB9B8

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/06/2023
Supervised By :Jagrut Upadhyay 04/06/2023

Quant Time: Apr 06 00:58:07 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 06 00:55:13 2023
Response via : Initial Calibration



TIC: BP014561.D\data.ms

(4) Pyridine-d5 (S)

3.799min (-0.000) 0.05 ng/u1

response

414

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	0.52#
54.00	32.60	0.46#
0.00	0.00	0.00

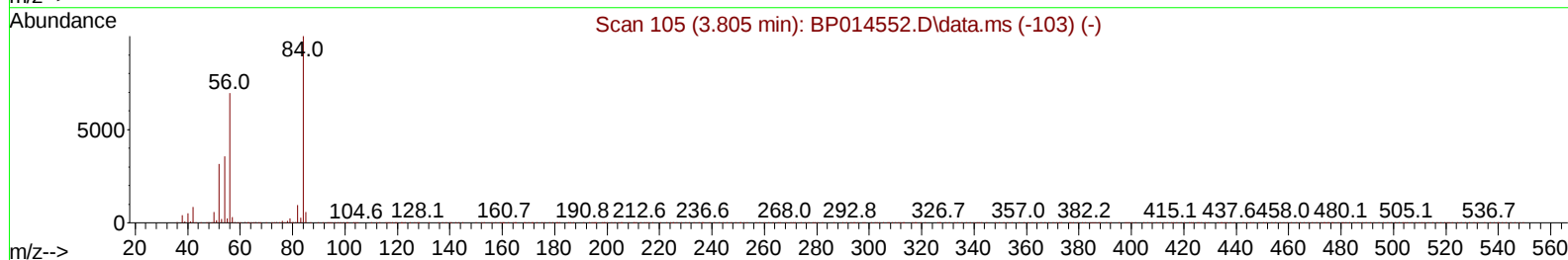
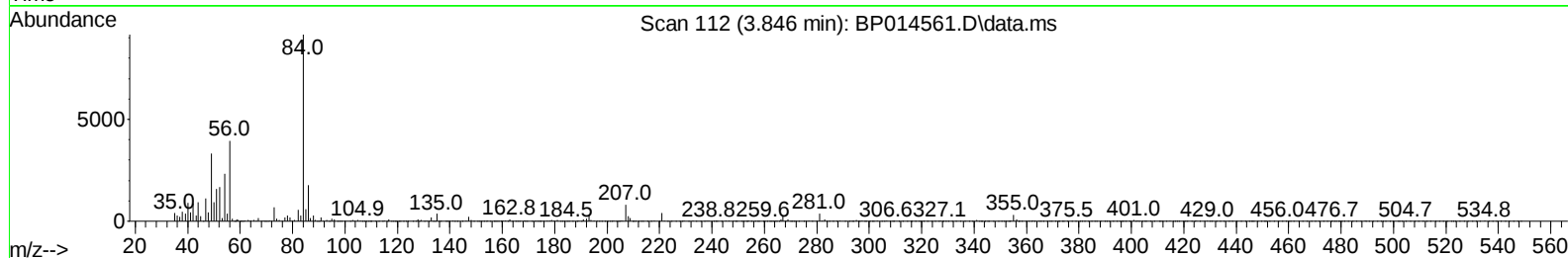
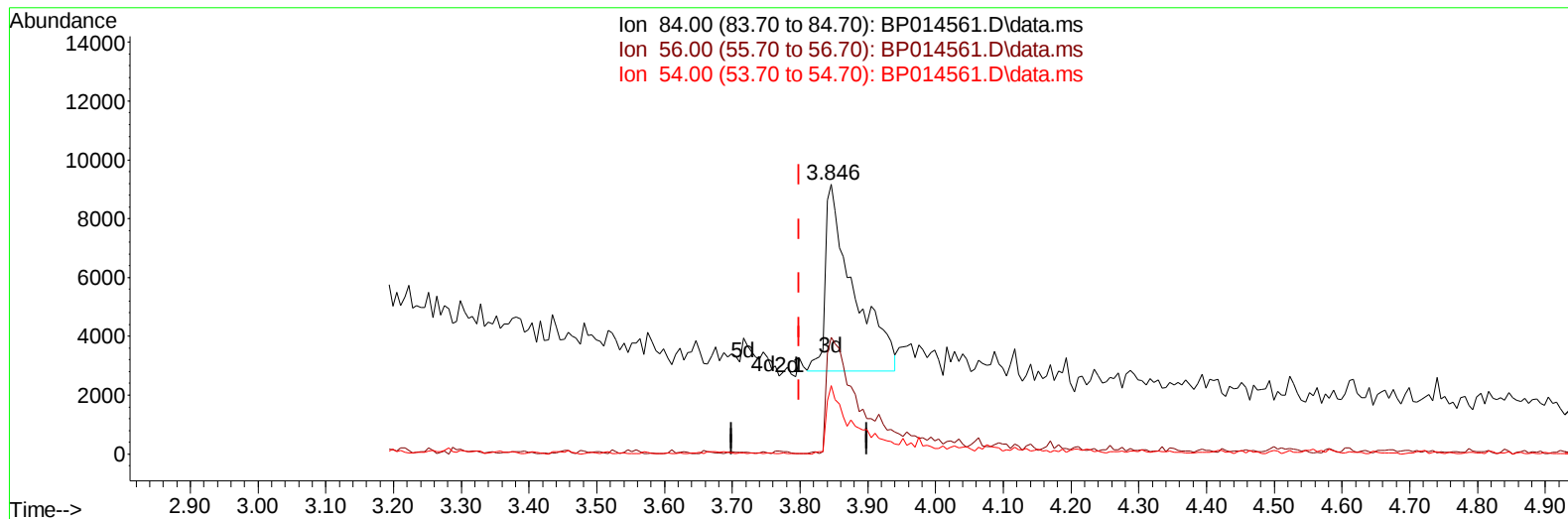
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TIC: BP014561.D\data.ms

(4) Pyridine-d5 (S)

3.846min (+ 0.047) 2.40 ng/ul m

response 18302

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	43.19#
54.00	32.60	25.32#
0.00	0.00	0.00

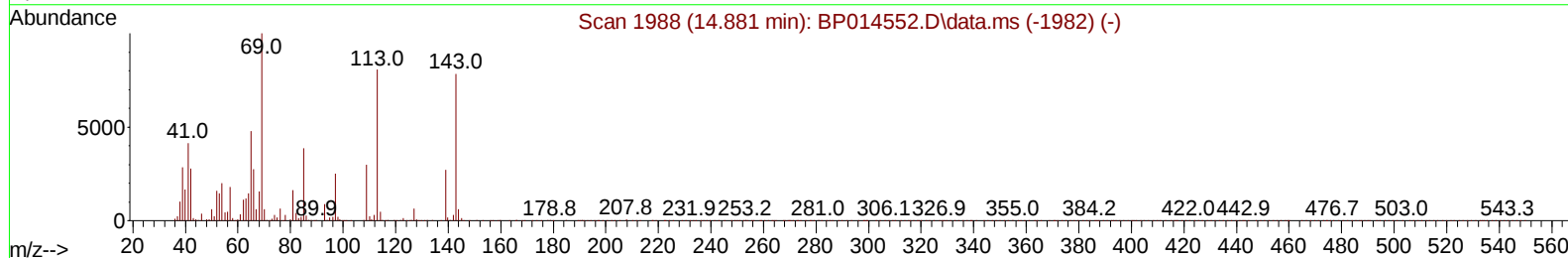
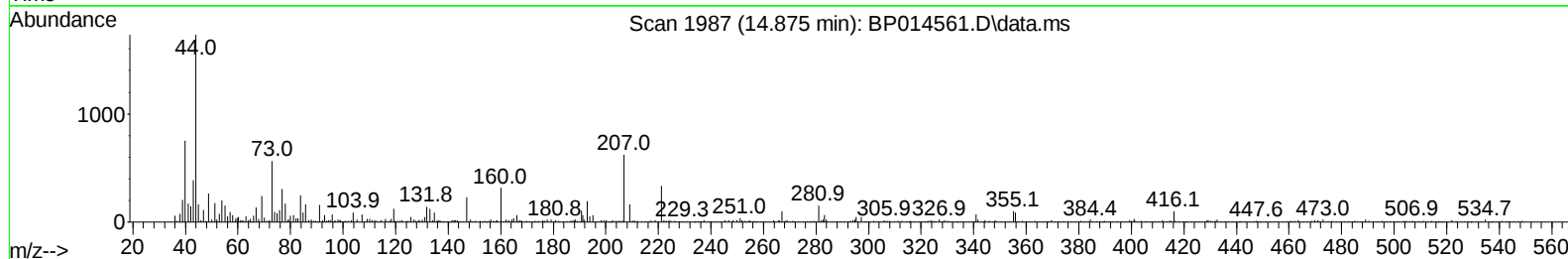
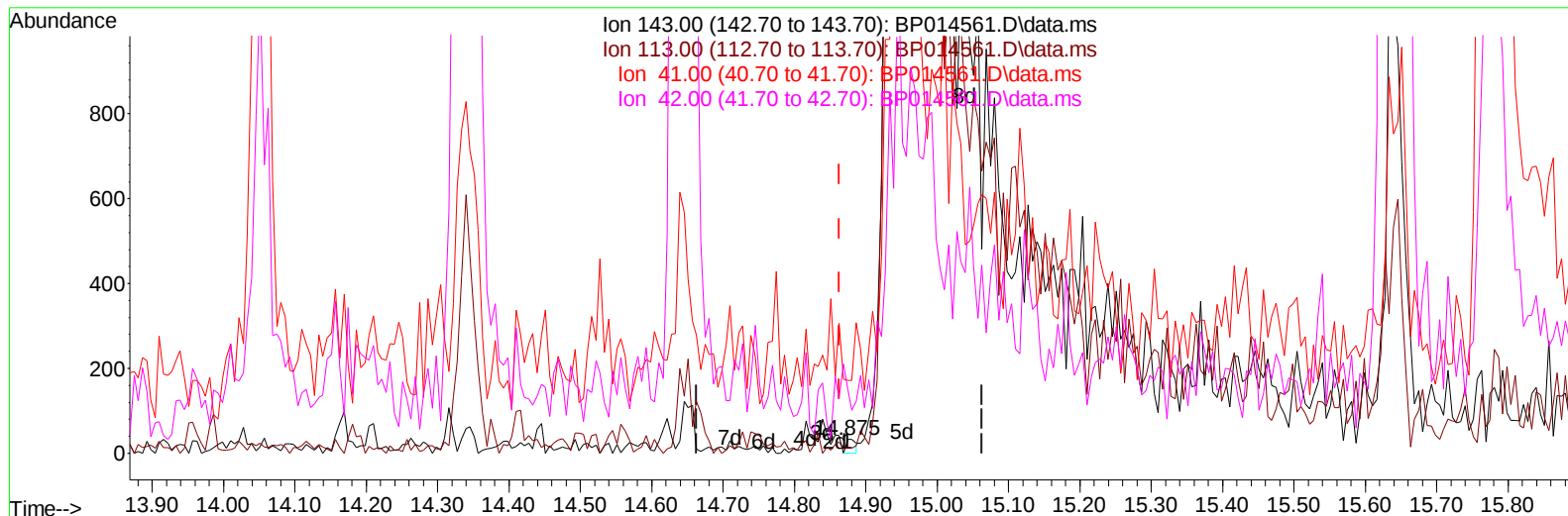
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
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TIC: BP014561.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.875min (+ 0.012) 0.01 ng/u1

response 31

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	42.42#
41.00	44.20	521.21#
42.00	31.50	451.52#

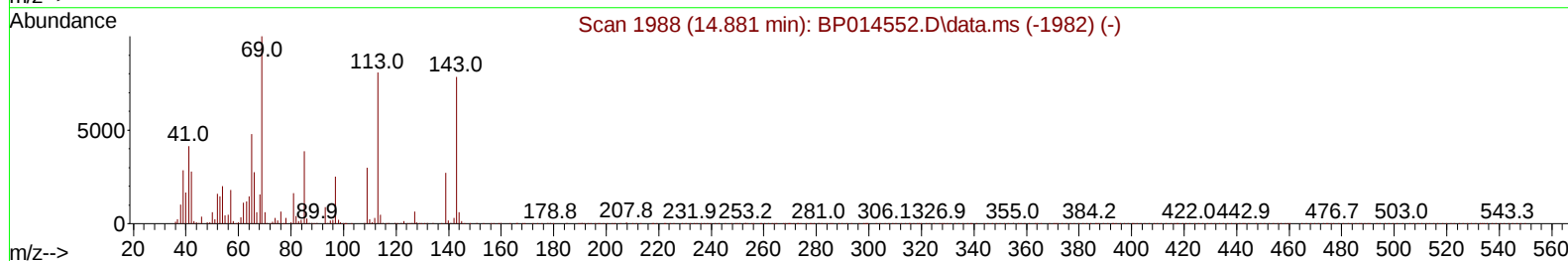
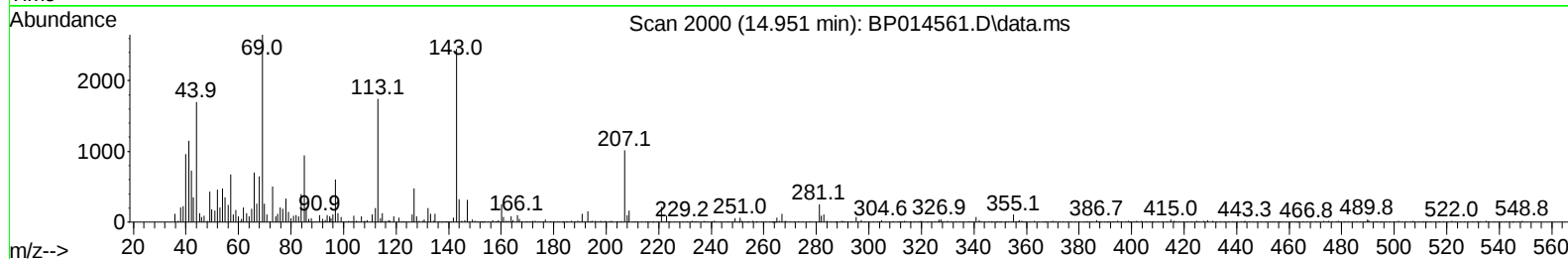
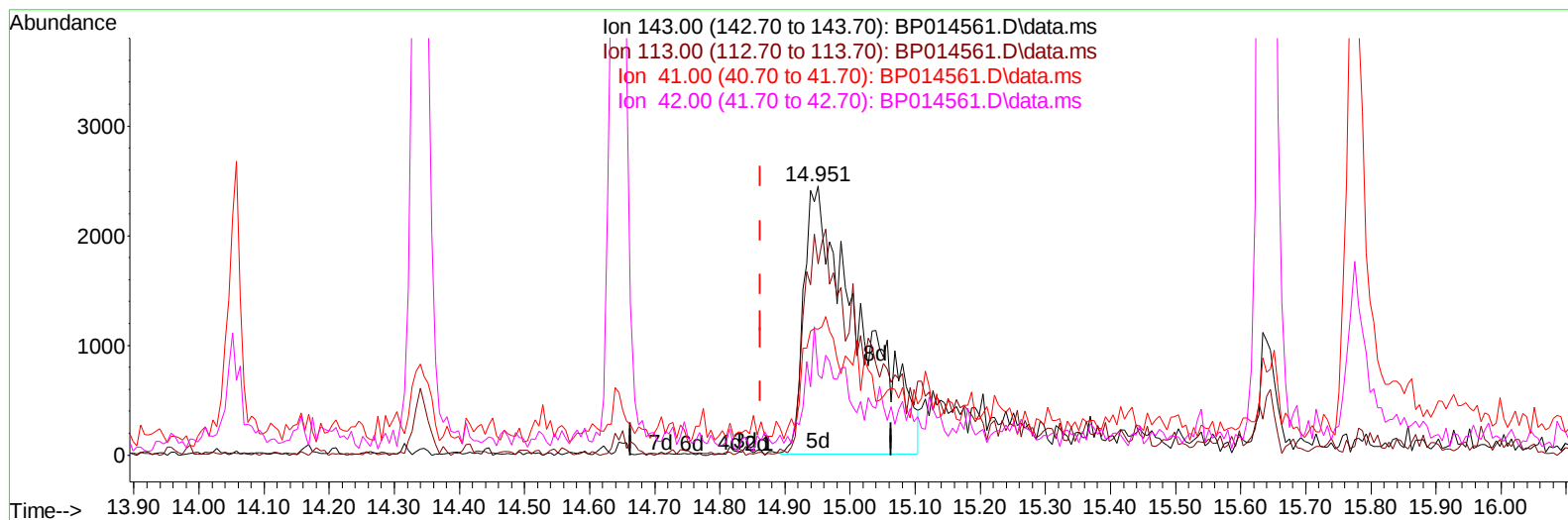
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Instrument :
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TIC: BP014561.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.951min (+ 0.088) 3.67 ng/ul m

response 14395

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	71.17#
41.00	44.20	46.82
42.00	31.50	29.85

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.993	152	115364	20.000	ng/ul	0.00
20) Naphthalene-d8	10.816	136	490304	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.645	164	320404	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.404	188	686385	20.000	ng/ul	0.00
79) Chrysene-d12	21.492	240	654537	20.000	ng/ul	0.00
88) Perylene-d12	23.998	264	747772	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	13886	4.668	ng/uL	0.00
4) Pyridine-d5	3.846	84	18302m	2.401	ng/ul	0.05
7) Phenol-d5	7.175	99	52665	4.950	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.322	67	206369	30.231	ng/ul	0.01
11) 2-Chlorophenol-d4	7.528	132	171679	22.325	ng/ul	0.01
15) 4-Methylphenol-d8	8.722	113	104656	12.195	ng/ul	0.01
21) Nitrobenzene-d5	9.175	128	111854	28.259	ng/ul	0.01
24) 2-Nitrophenol-d4	9.899	143	114859	25.342	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.451	165	187296	22.691	ng/ul	0.02
31) 4-Chloroaniline-d4	10.975	131	113681	10.456	ng/ul	0.02
46) Dimethylphthalate-d6	14.057	166	869325	33.463	ng/ul	0.01
49) Acenaphthylene-d8	14.339	160	913882	31.703	ng/ul	0.00
54) 4-Nitrophenol-d4	14.951	143	14395m	3.668	ng/ul	0.09
60) Fluorene-d10	15.639	176	706462	32.609	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	97295	19.264	ng/ul	0.01
73) Anthracene-d10	17.504	188	1169520	35.207	ng/ul	0.00
81) Pyrene-d10	19.733	212	1387418	31.557	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264	1378695	34.703	ng/ul	0.00

Target Compounds Qvalue

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

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