

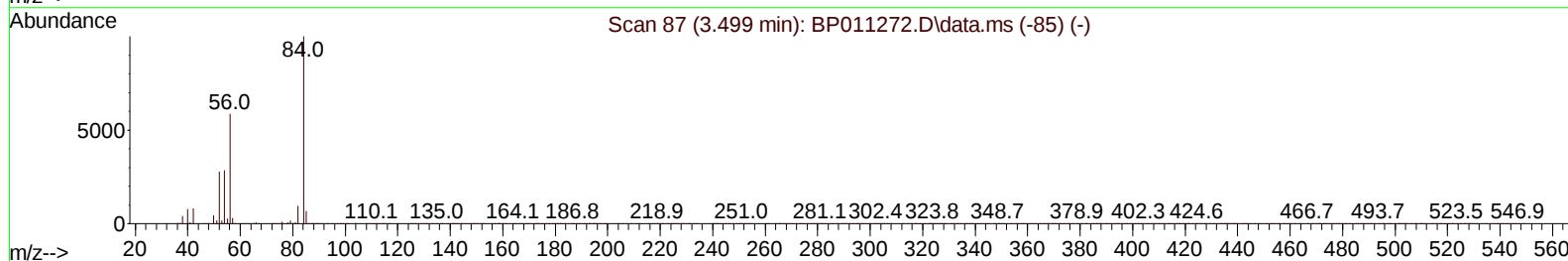
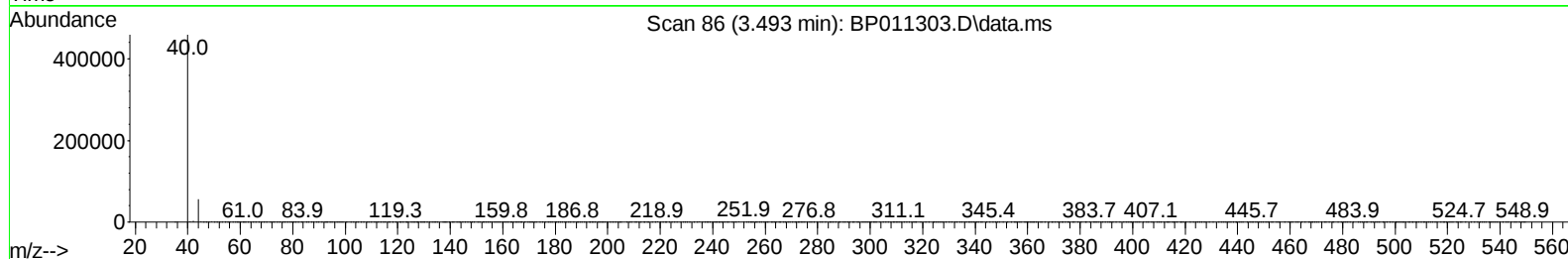
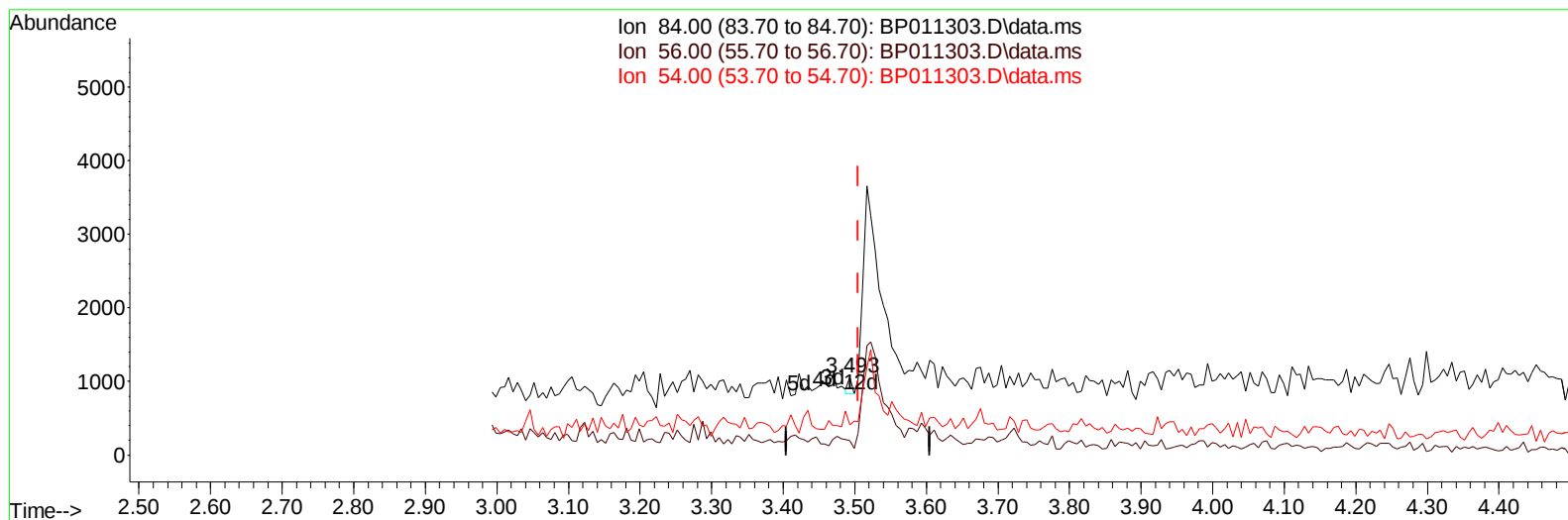
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
 Data File : BP011303.D
 Acq On : 06 Aug 2022 06:40
 Operator : CG/JU
 Sample : N4026-22
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJJ7

Manual IntegrationsAPPROVED

Quant Time: Aug 08 01:09:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080222.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 02 23:52:37 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/08/2022
 Supervised By :Sohil Jodhani 08/10/2022



TIC: BP011303.D\data.ms

(4) Pyridine-d5 (S)

3.493min (-0.012) 0.02 ng/u1

response 80

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.60	18.56#
54.00	33.20	38.92
0.00	0.00	0.00

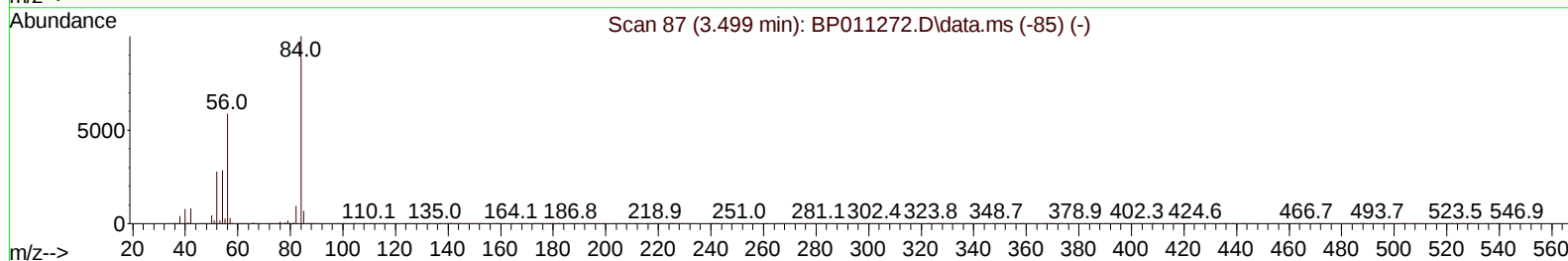
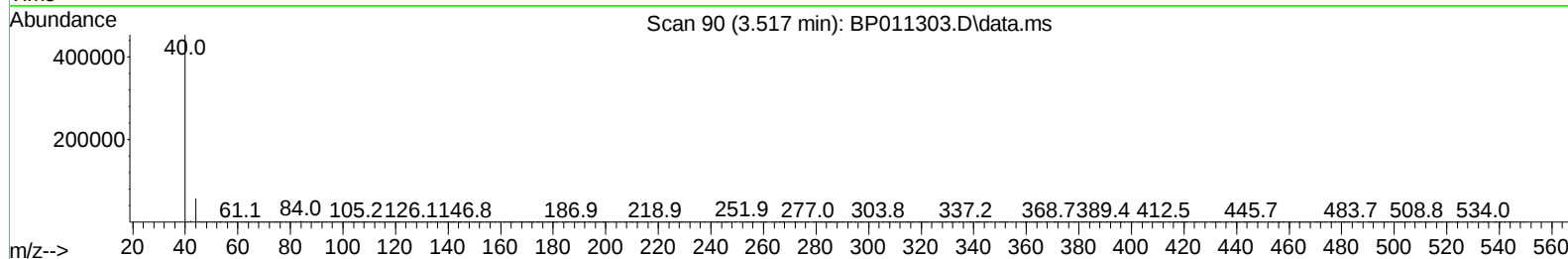
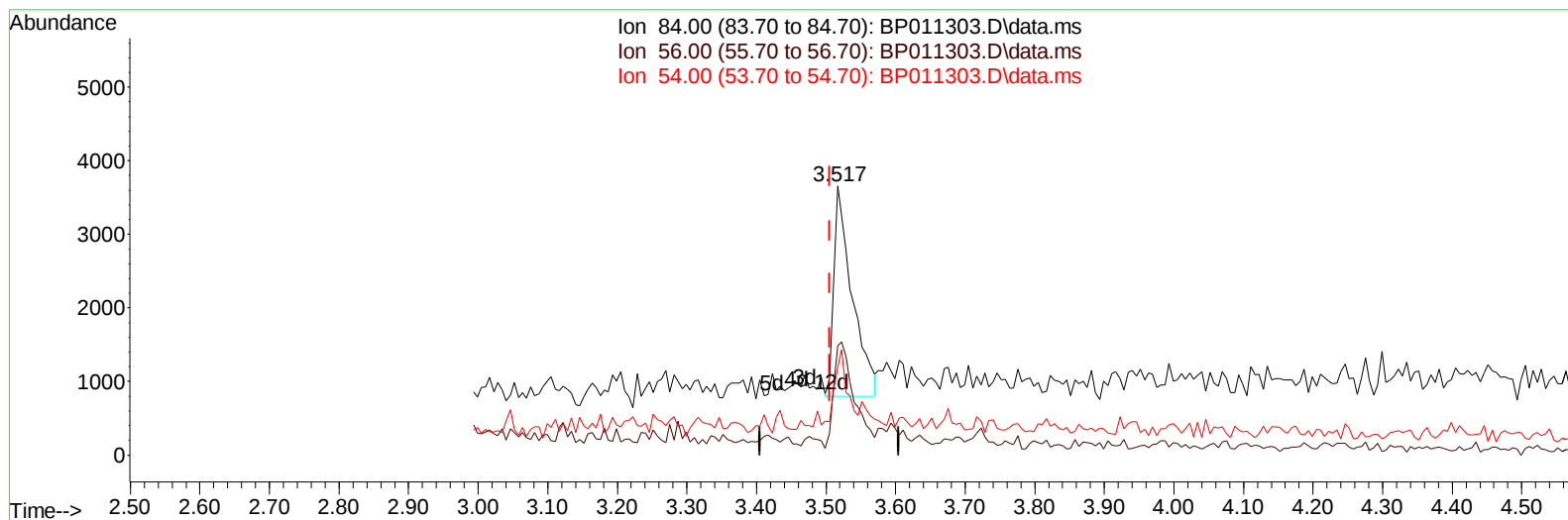
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(4) Pyridine-d5 (S)

3.517min (+ 0.012) 1.59 ng/ul m

response 5168

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.60	40.77#
54.00	33.20	32.61
0.00	0.00	0.00

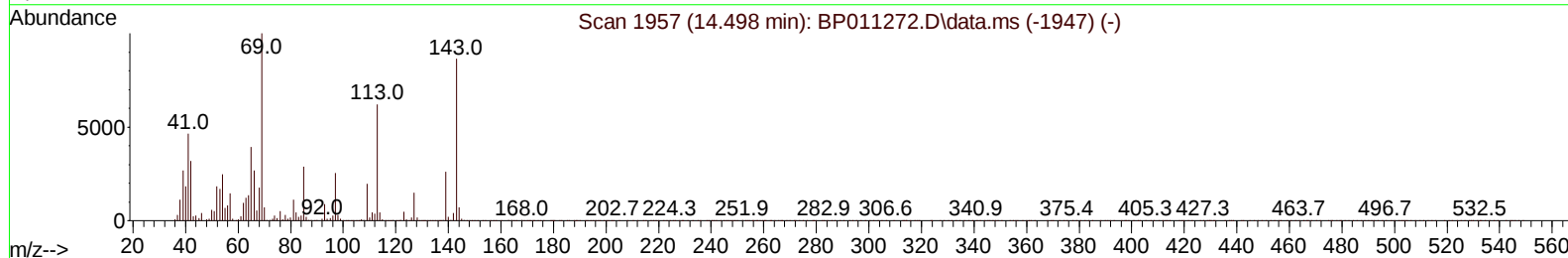
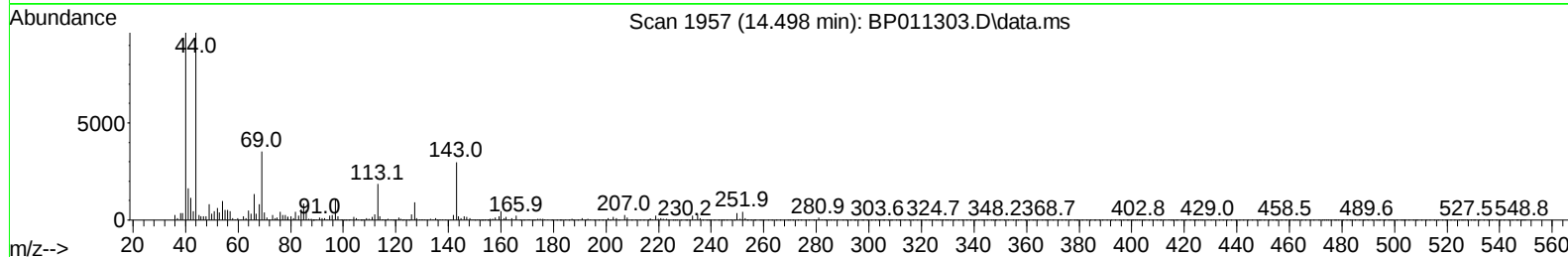
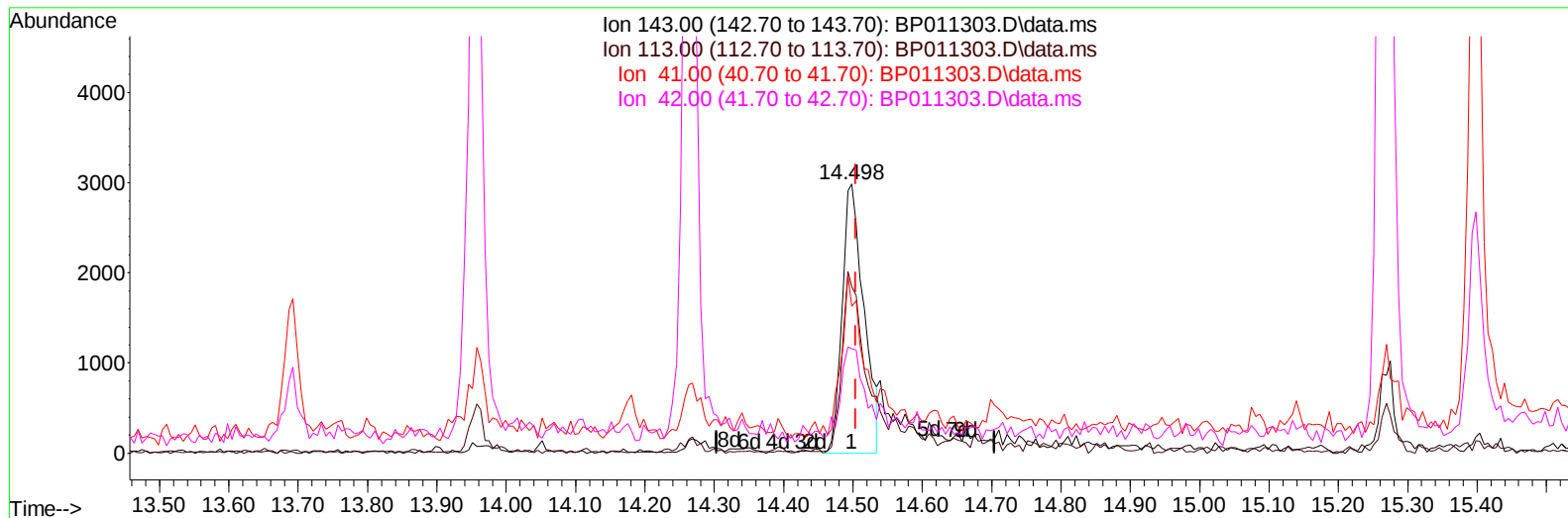
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TIC: BP011303.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.006) 2.00 ng/u1

response 6451

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	61.96
41.00	54.40	54.62
42.00	35.50	38.81

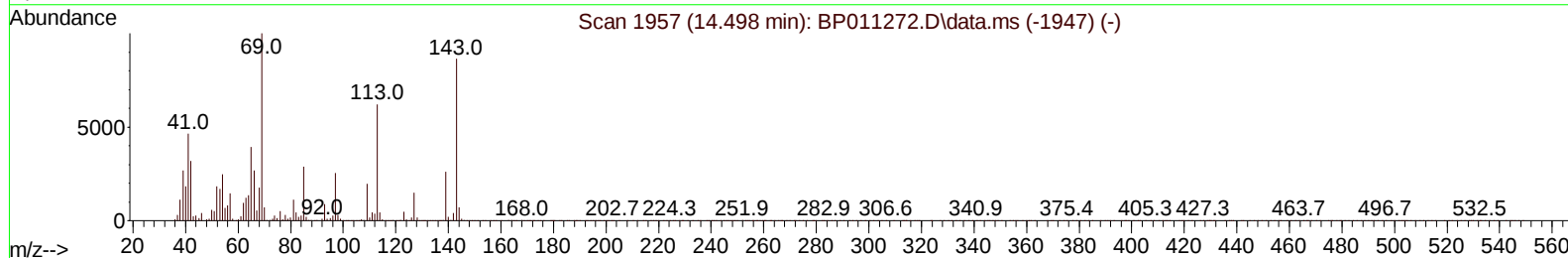
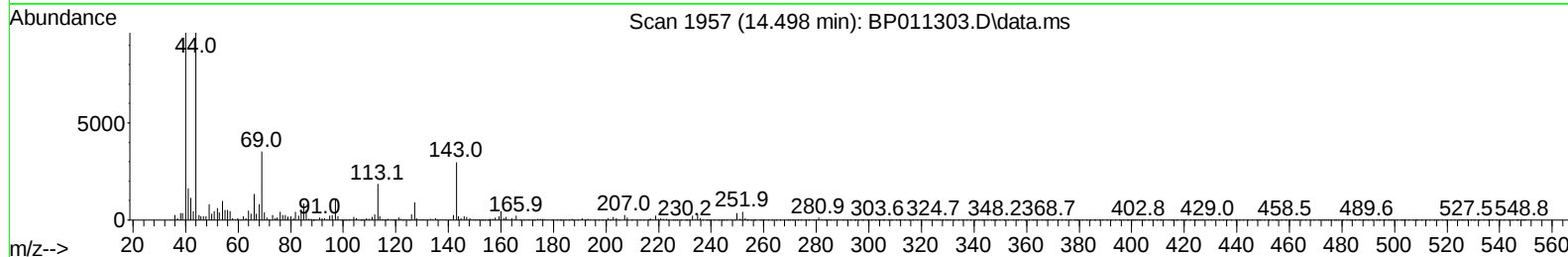
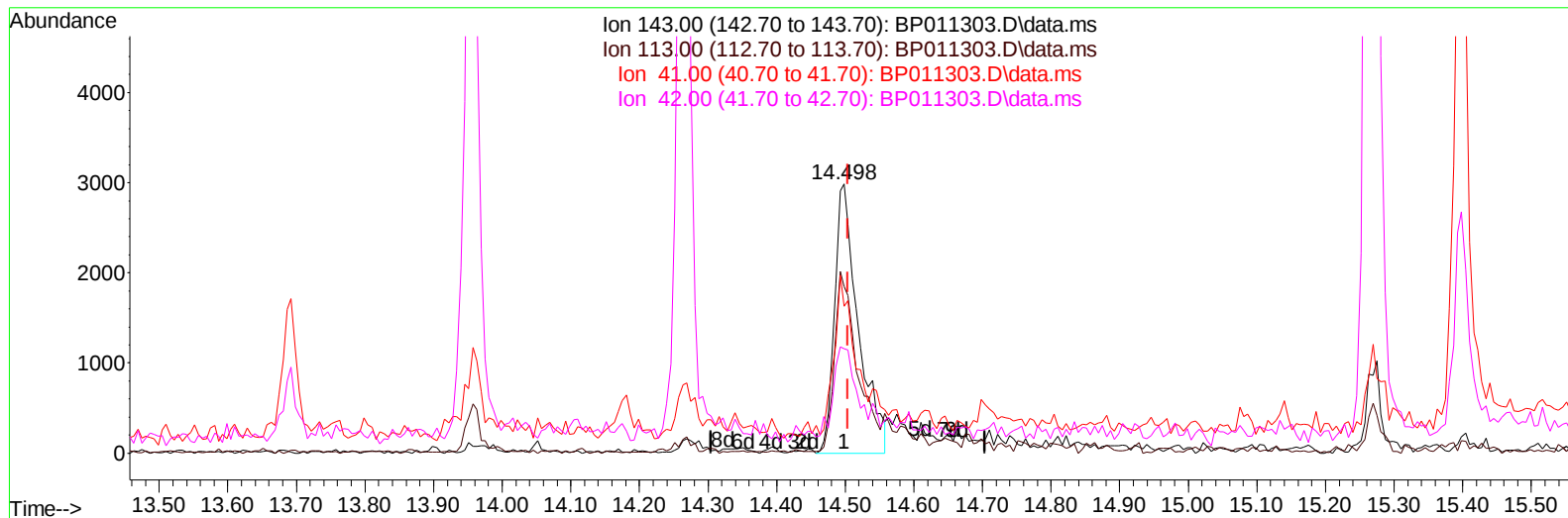
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Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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TIC: BP011303.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.006) 2.23 ng/ul m

response 7169

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	61.96
41.00	54.40	54.62
42.00	35.50	38.81

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	97473	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.387	136	420083	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.263	164	276450	20.000	ng/ul	#-0.01
64) Phenanthrene-d10	17.034	188	628641	20.000	ng/ul	-0.01
79) Chrysene-d12	21.157	240	757202	20.000	ng/ul	-0.02
88) Perylene-d12	23.468	264	730045	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	2461	2.041	ng/uL	0.00
4) Pyridine-d5	3.517	84	5168m	1.589	ng/ul	0.01
7) Phenol-d5	6.775	99	23100	2.844	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.940	67	156852	26.760	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.128	132	97081	14.662	ng/ul	-0.01
15) 4-Methylphenol-d8	8.310	113	48194	7.211	ng/ul	-0.02
21) Nitrobenzene-d5	8.758	128	123306	40.809	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.475	143	76514	27.015	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.010	165	110729	20.502	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.540	131	49424	5.915	ng/ul	0.00
46) Dimethylphthalate-d6	13.693	166	460517	24.322	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	743507	31.210	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.498	143	7169m	2.227	ng/ul	0.00
60) Fluorene-d10	15.269	176	662241	41.105	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.398	200	67033	22.964	ng/ul	-0.02
73) Anthracene-d10	17.134	188	1195376	43.266	ng/ul	-0.01
81) Pyrene-d10	19.392	212	1399073	36.544	ng/ul	-0.02
92) Benzo(a)pyrene-d12	23.321	264	1224874	33.670	ng/ul	-0.02

Target Compounds Qvalue

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

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