

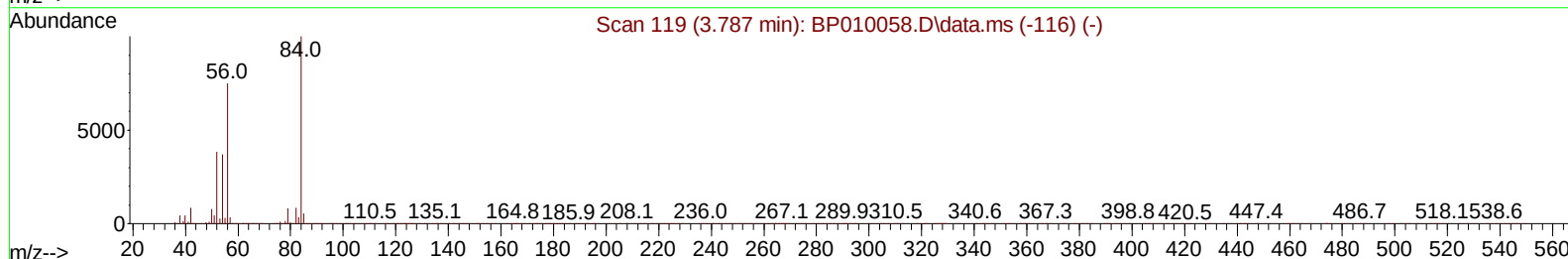
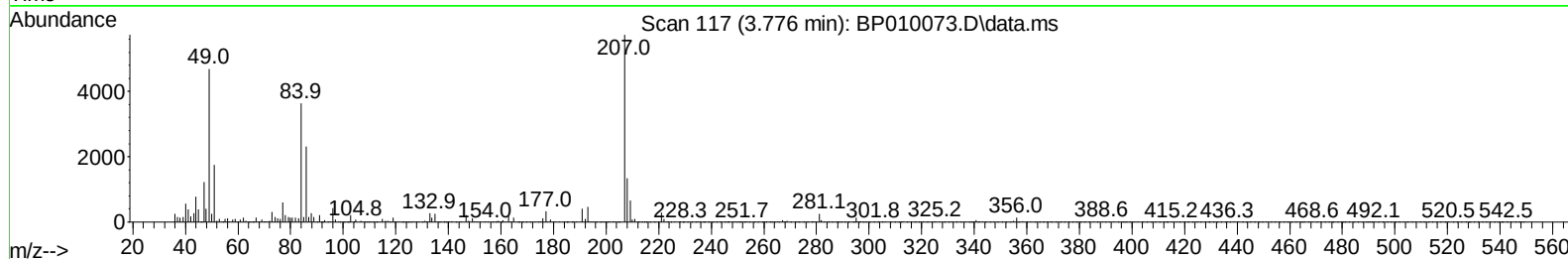
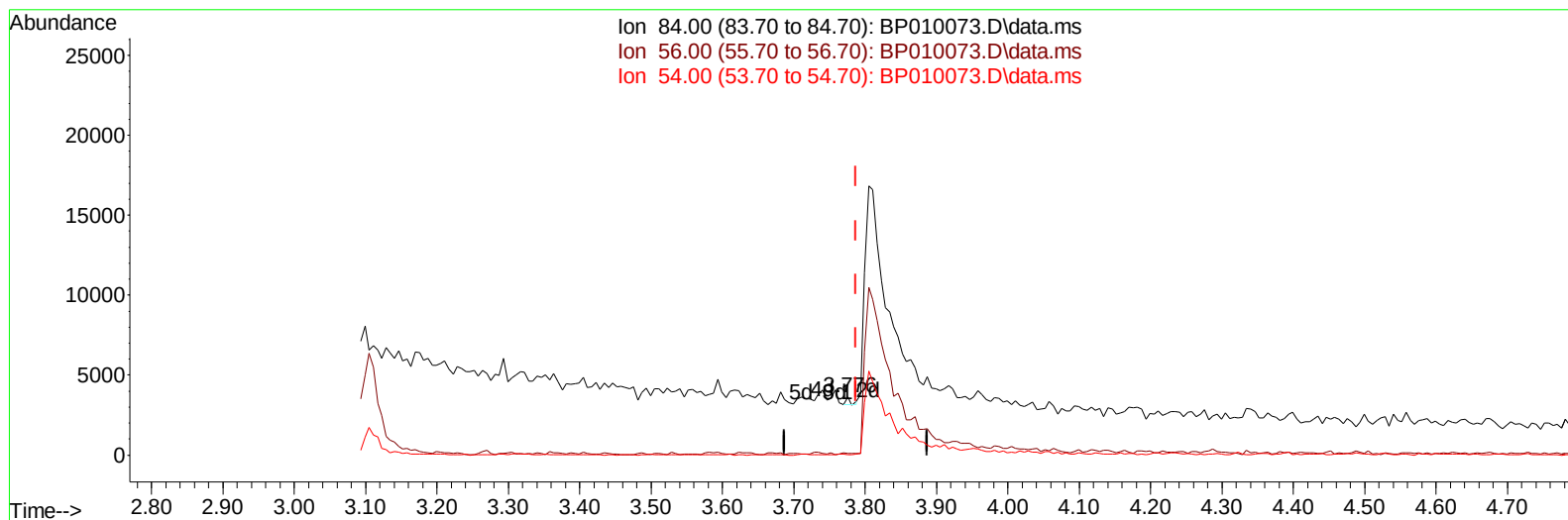
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010073.D
 Acq On : 23 Apr 2022 01:36
 Operator : CG/JU
 Sample : N2475-08
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A0BN8

Manual IntegrationsAPPROVED

Quant Time: Apr 23 02:25:31 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 14:49:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/25/2022
 Supervised By :mohammad ahmed 04/26/2022



TIC: BP010073.D\data.ms

(4) Pyridine-d5 (S)

3.776min (-0.011) 0.02 ng/ul

response 212

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	3.63#
54.00	24.80	0.63#
0.00	0.00	0.00

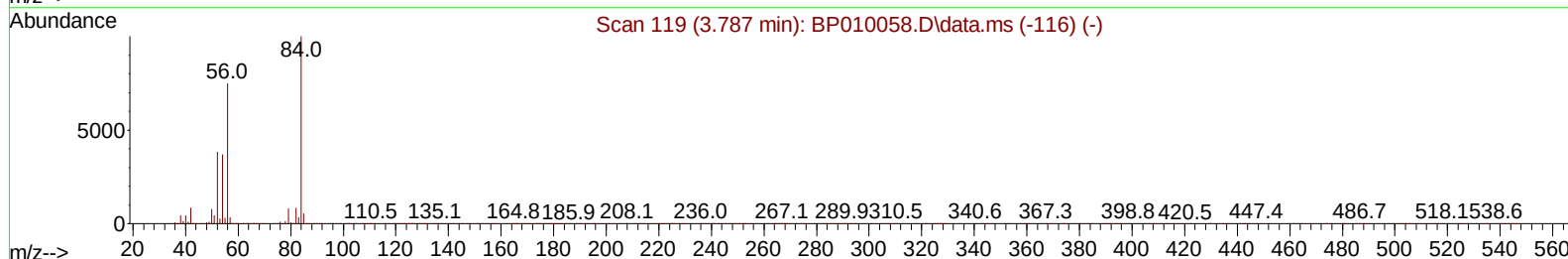
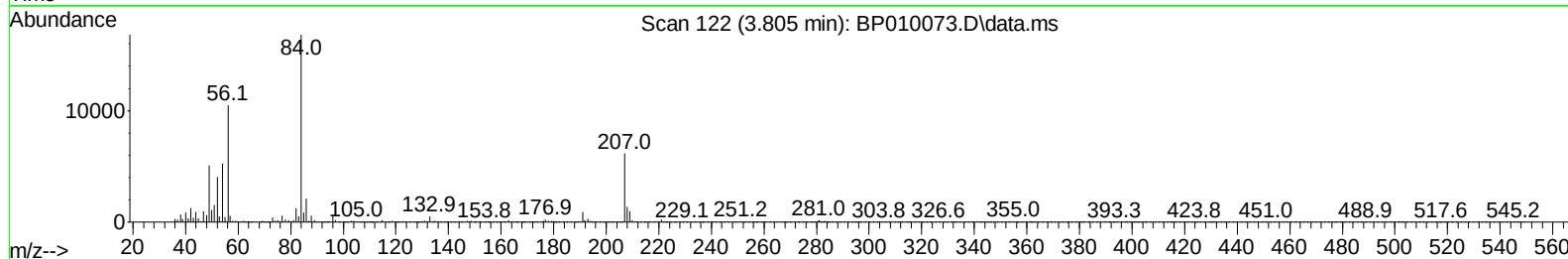
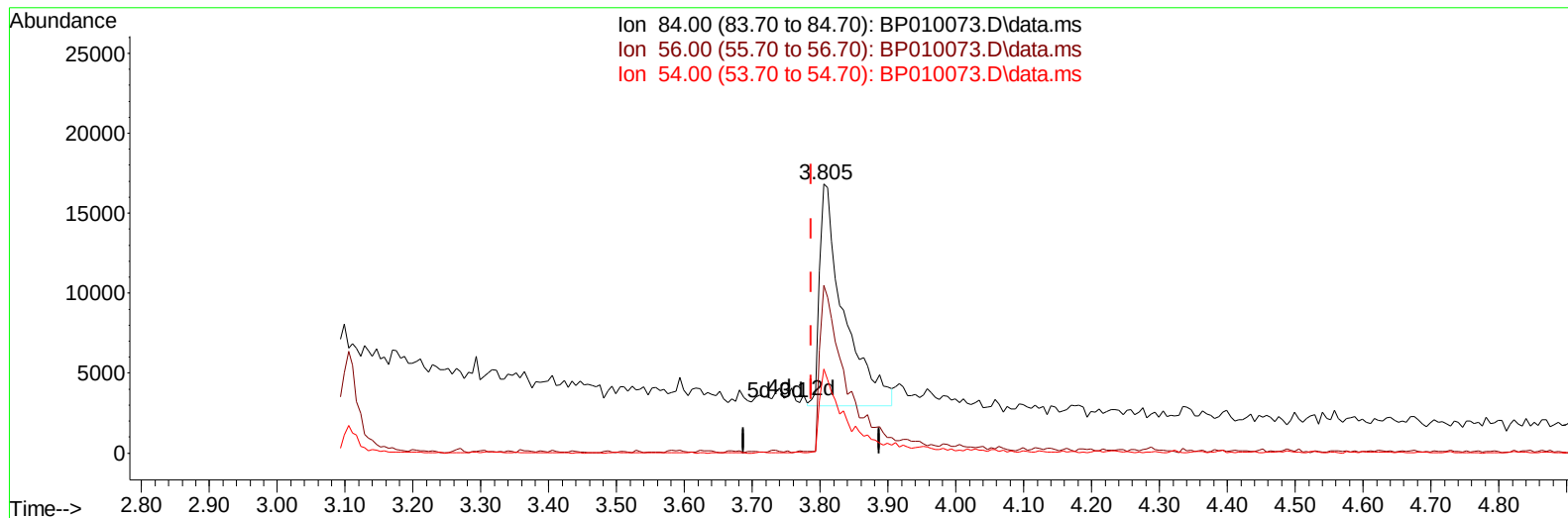
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(4) Pyridine-d5 (S)

3.805min (+ 0.018) 3.28 ng/ul m

response 34524

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	62.38#
54.00	24.80	31.16#
0.00	0.00	0.00

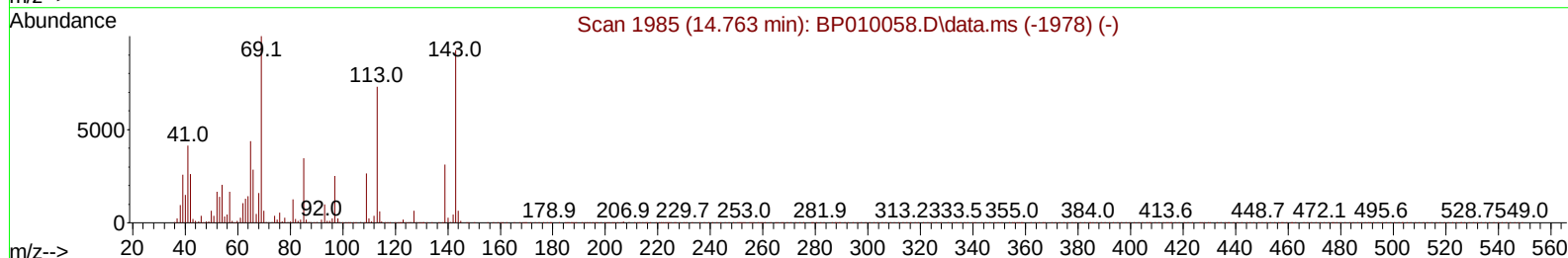
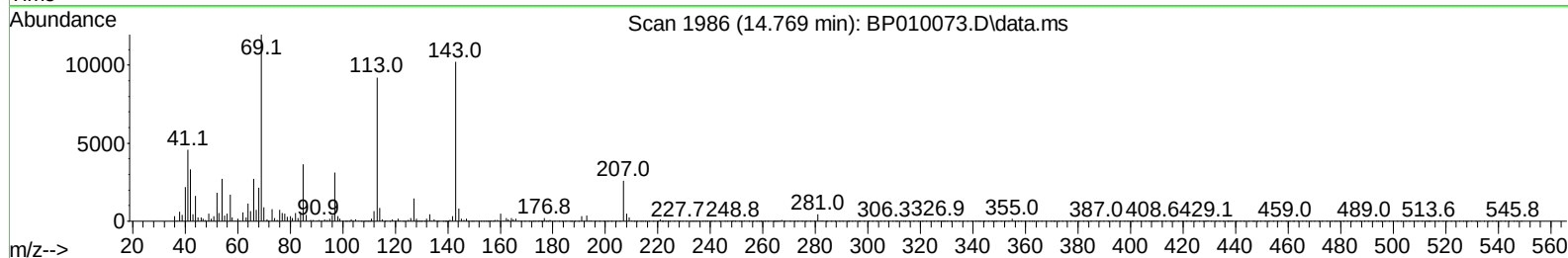
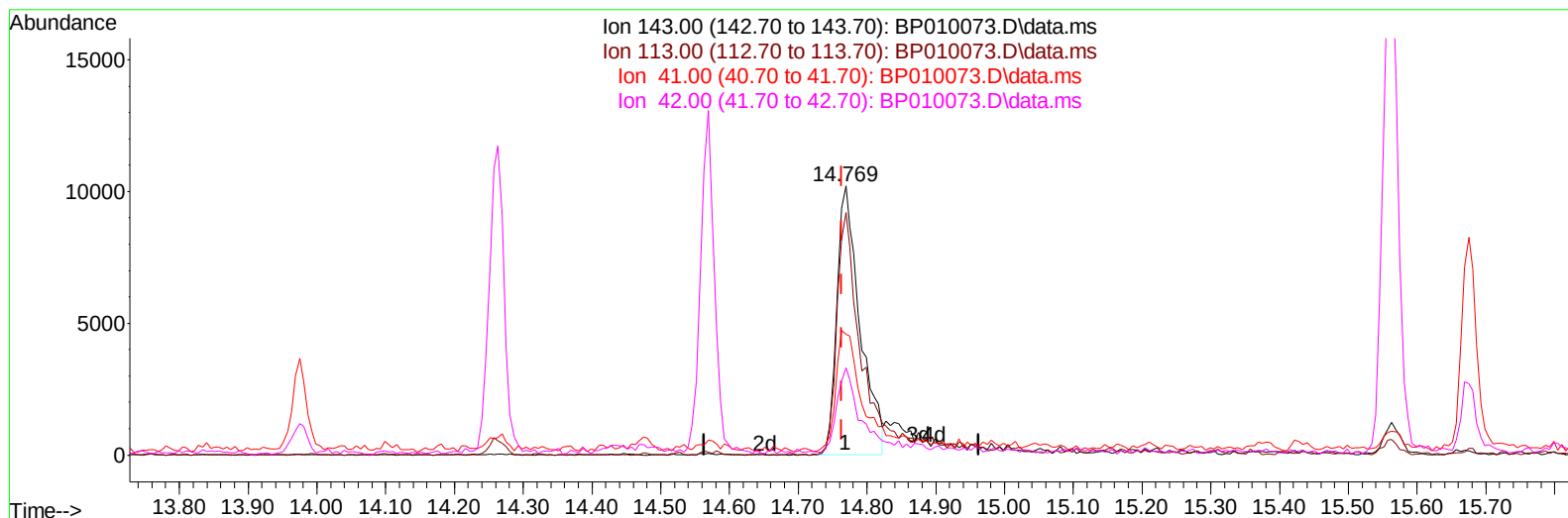
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ALS Vial : 69 Sample Multiplier: 1

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TIC: BP010073.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (+ 0.006) 3.05 ng/ul

response 23848

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	90.13#
41.00	23.20	44.84#
42.00	18.00	32.53#

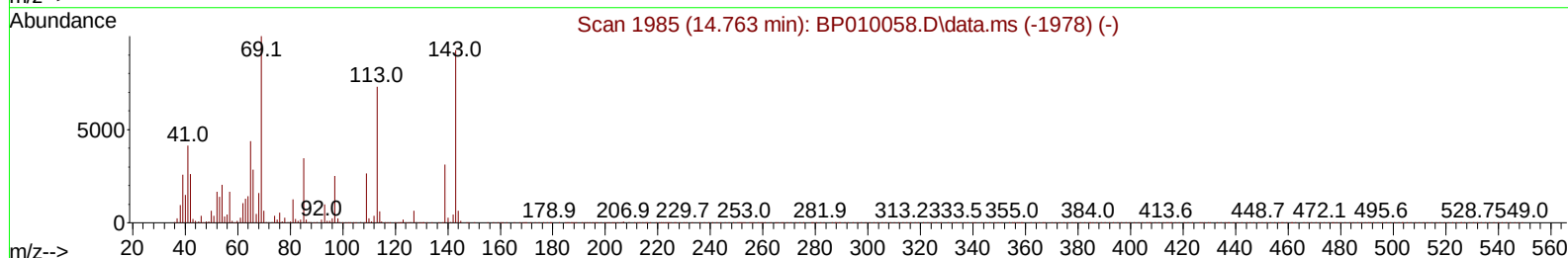
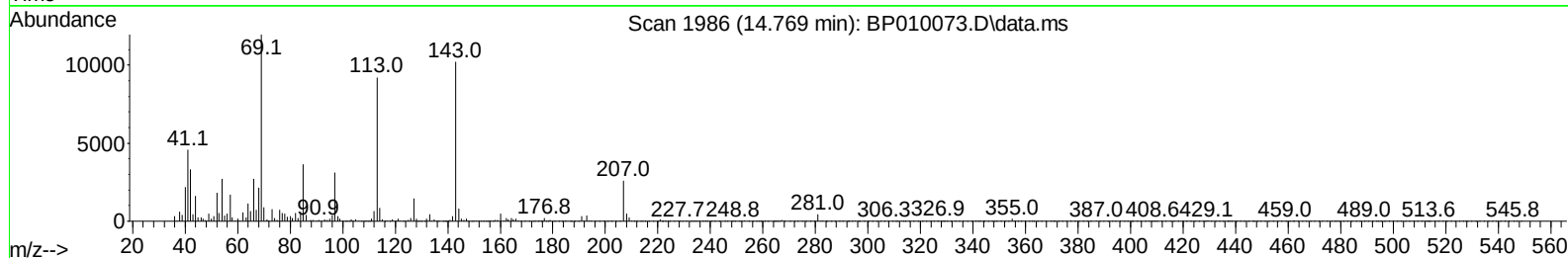
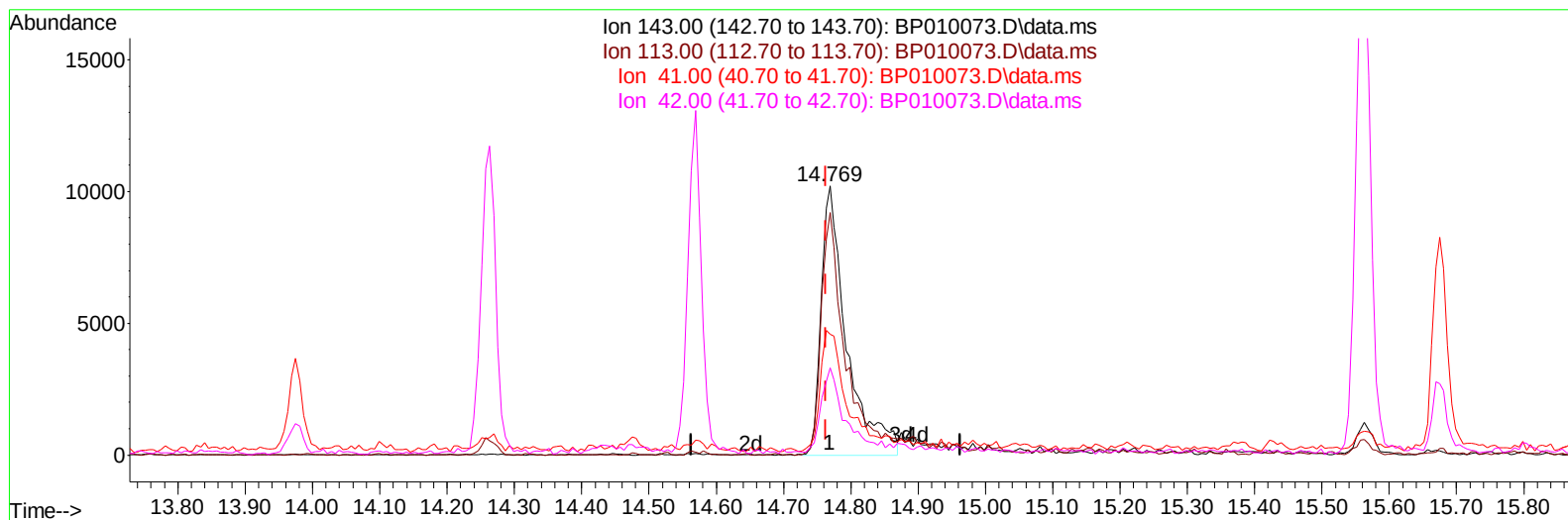
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Misc :
ALS Vial : 69 Sample Multiplier: 1

Instrument :
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TIC: BP010073.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (+ 0.006) 3.43 ng/ul m

response 26798

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	90.13#
41.00	23.20	44.84#
42.00	18.00	32.53#

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 ALS Vial : 69 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	163214	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	767559	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	545581	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	1137022	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	852377	20.000	ng/ul	# 0.00
88) Perylene-d12	23.857	264	757544	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	16628	4.443	ng/uL	0.00
4) Pyridine-d5	3.805	84	34524m	3.279	ng/ul	0.02
7) Phenol-d5	7.099	99	97599	6.574	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67	247630	27.798	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	247884	22.546	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	185197	15.056	ng/ul	-0.01
21) Nitrobenzene-d5	9.099	128	165564	27.121	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	171345	25.700	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	313035	24.118	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	236980	12.734	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	1293598	30.481	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1400505	27.266	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	26798m	3.430	ng/ul	0.00
60) Fluorene-d10	15.563	176	1117612	30.613	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	131780	18.329	ng/ul	0.00
73) Anthracene-d10	17.422	188	1748340	31.783	ng/ul	0.00
81) Pyrene-d10	19.651	212	1995962	41.175	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.692	264	1283704	32.323	ng/ul	0.00

Target Compounds Qvalue

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

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