

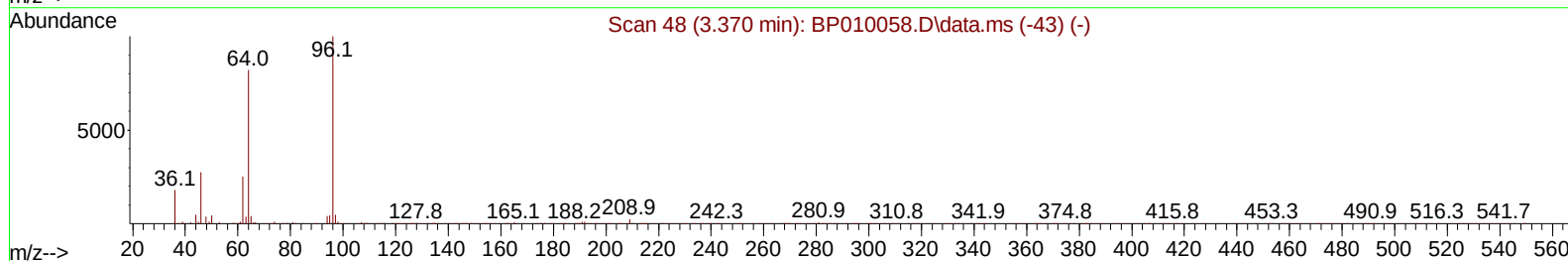
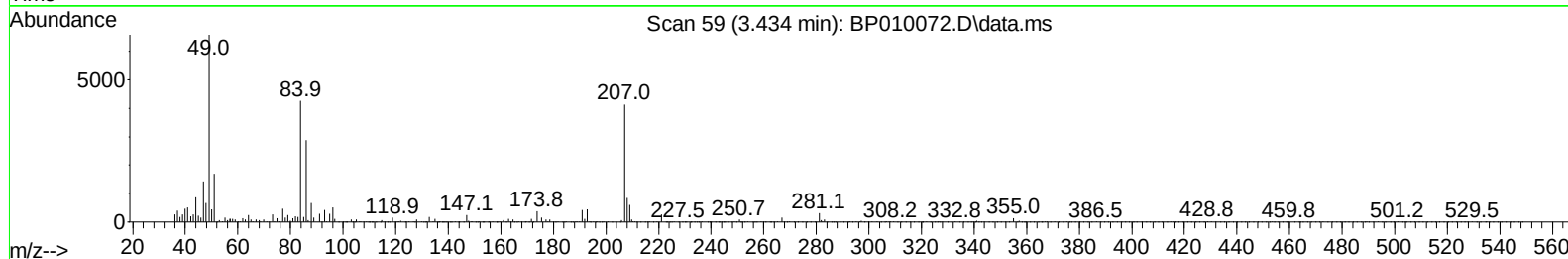
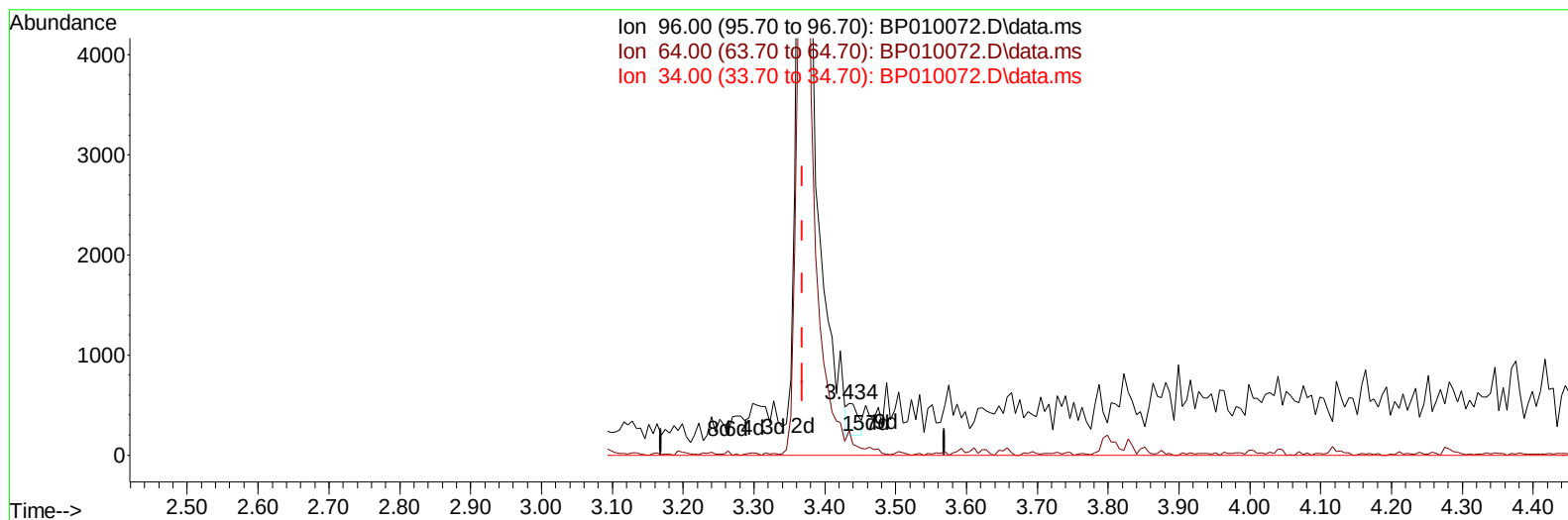
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010072.D
 Acq On : 23 Apr 2022 01:03
 Operator : CG/JU
 Sample : N2475-07
 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A0BM6

Manual IntegrationsAPPROVED

Quant Time: Apr 23 01:59:02 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: BP010072.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.434min (+ 0.065) 0.10 ng/uL

response 347

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	47.77
34.00	0.00	0.00
0.00	0.00	0.00

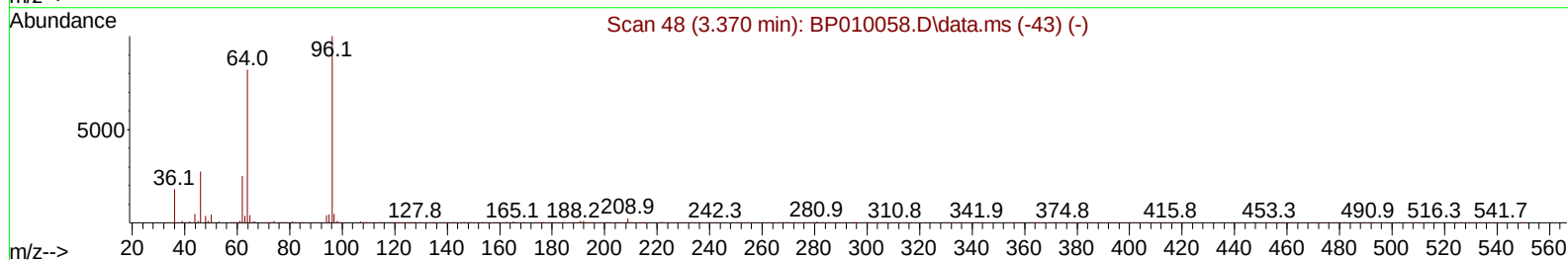
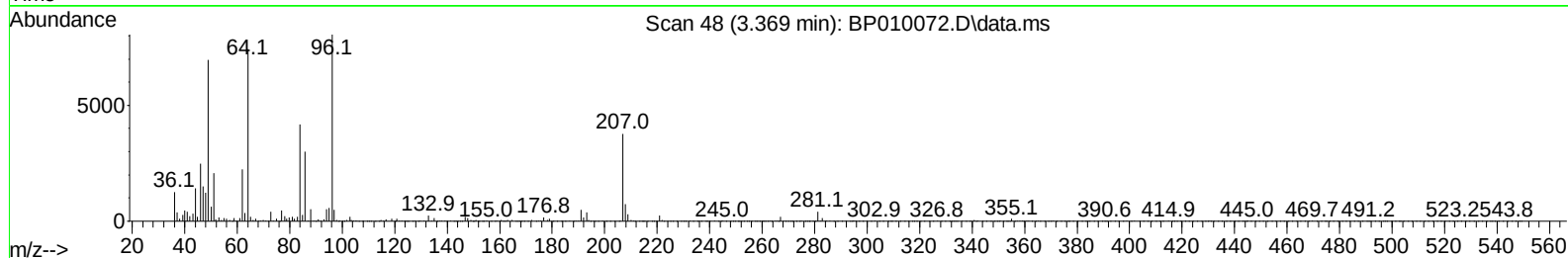
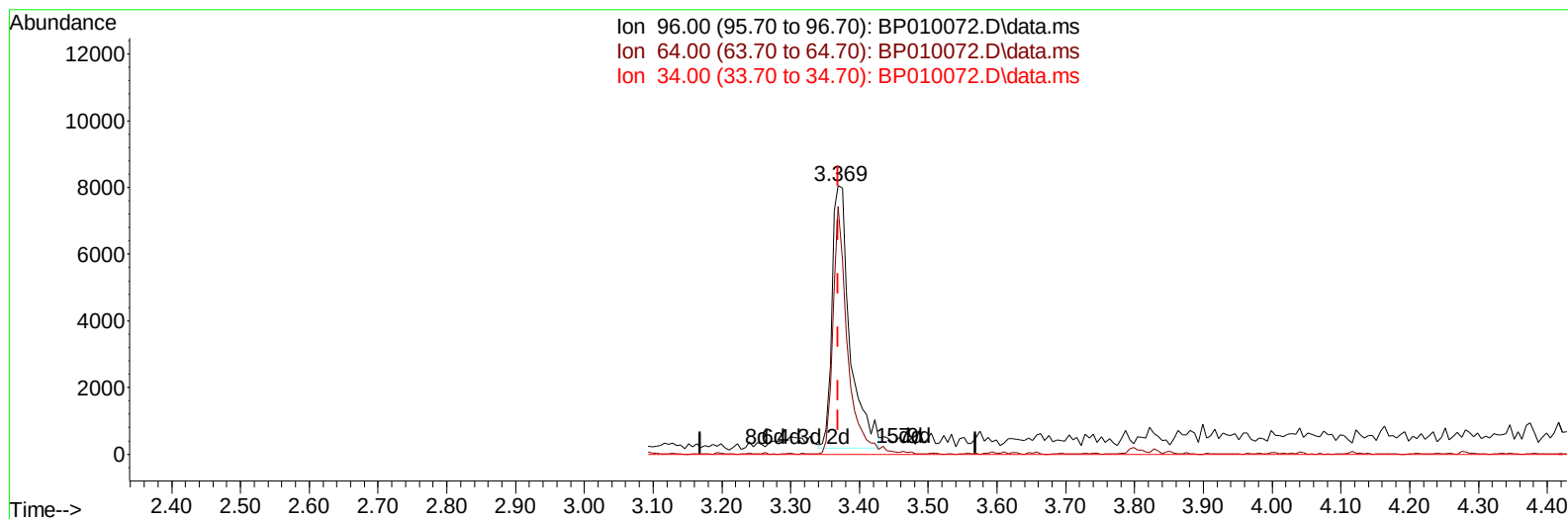
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(3) 1,4-Dioxane-d8 (S)

3.369min (+ 0.000) 4.17 ng/uL m

response 14218

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	92.34#
34.00	0.00	0.00
0.00	0.00	0.00

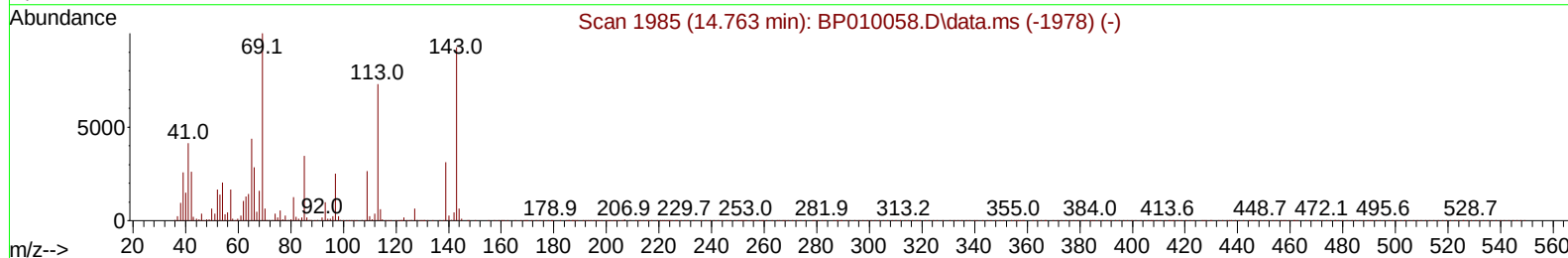
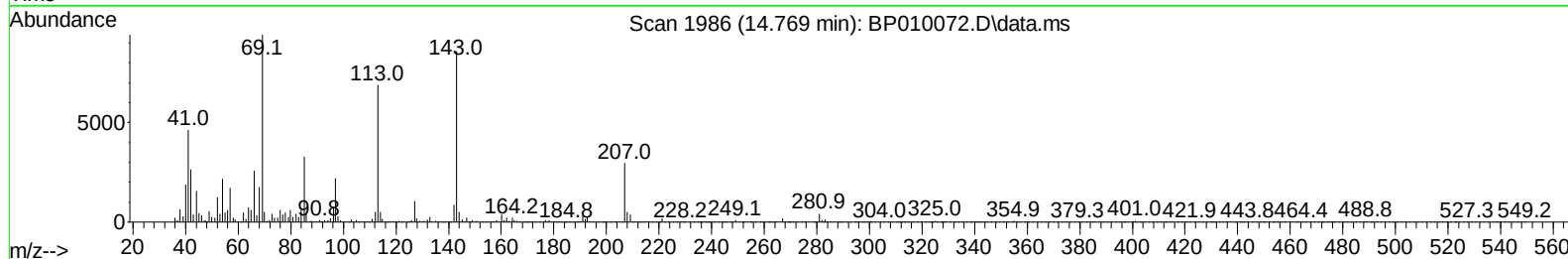
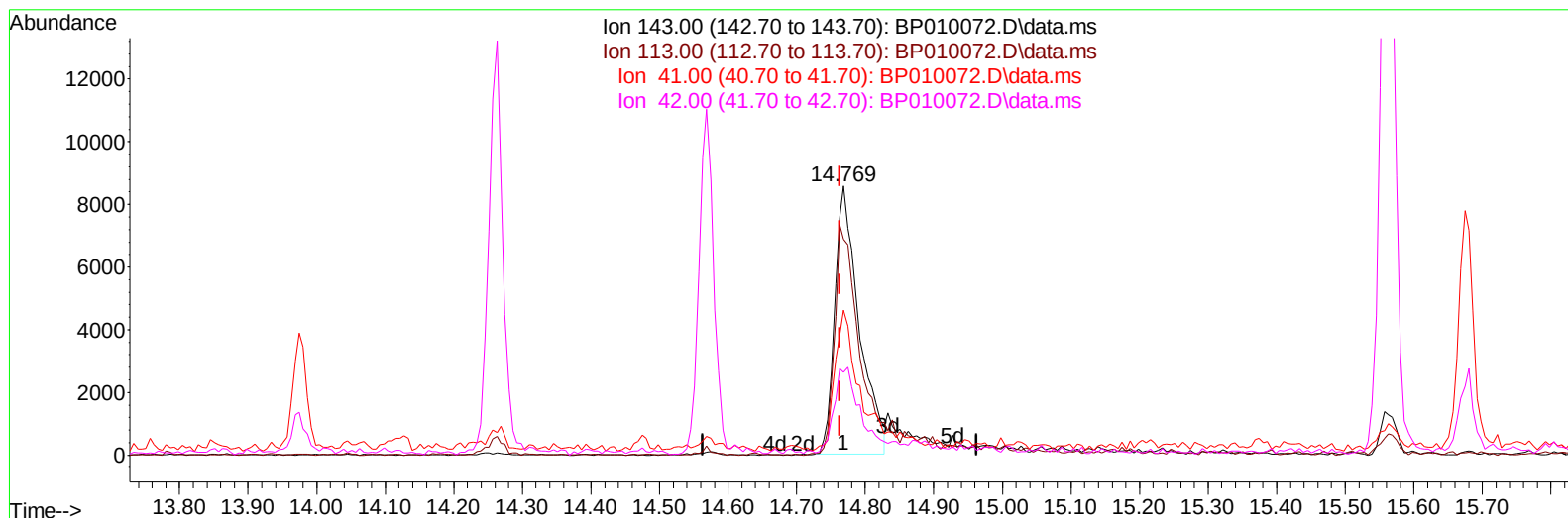
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Misc :
ALS Vial : 68 Sample Multiplier: 1

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TIC: BP010072.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (+ 0.006) 2.86 ng/ul

response 20967

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.24#
41.00	23.20	53.98#
42.00	18.00	30.76#

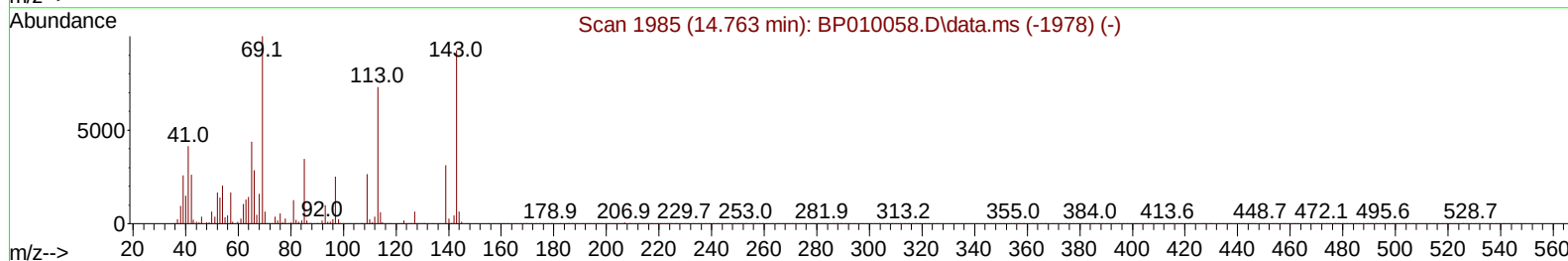
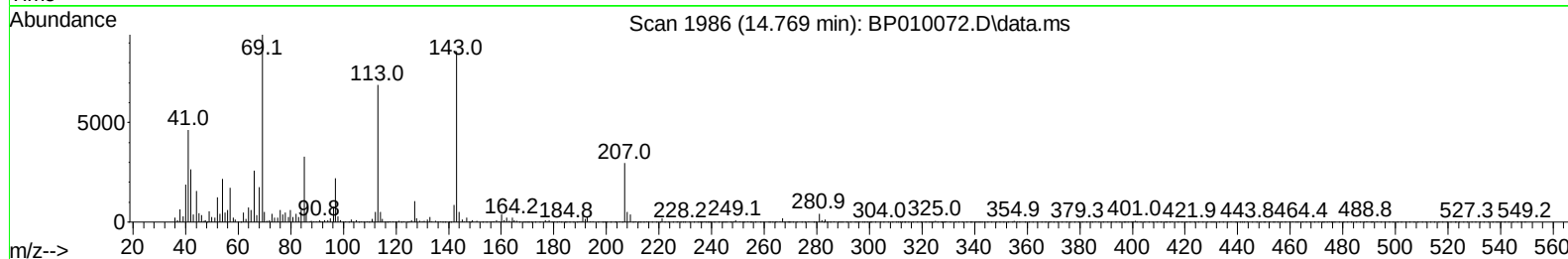
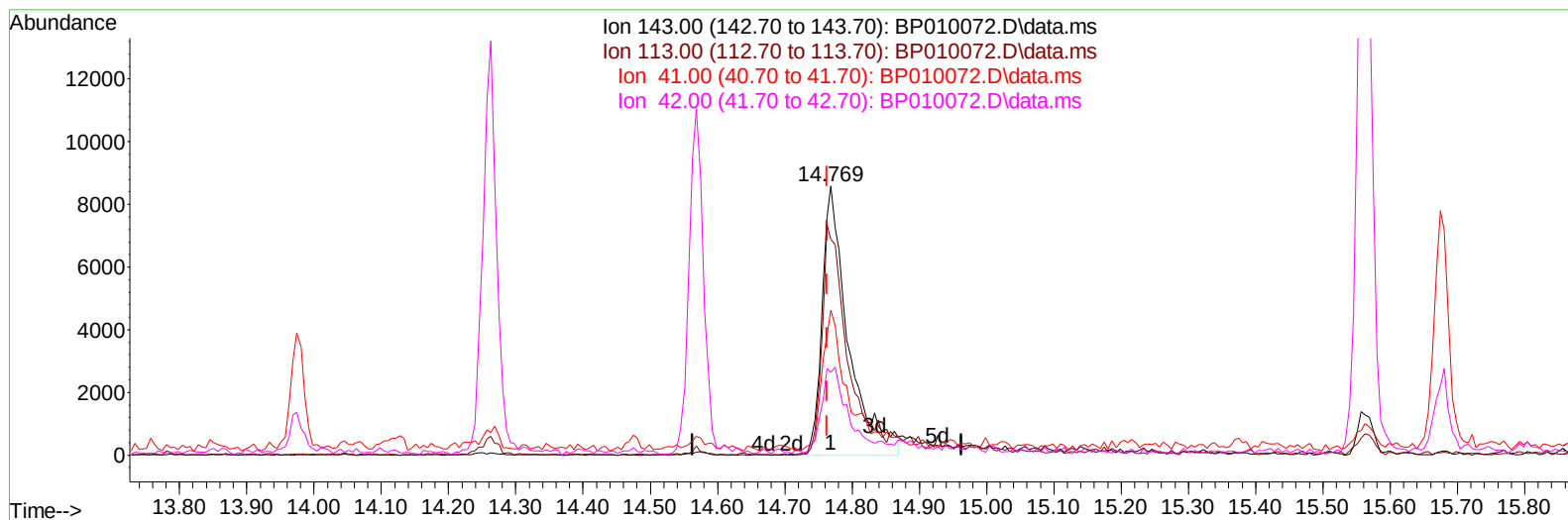
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 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
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TIC: BP010072.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (+ 0.006) 3.16 ng/ul m

response 23188

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.24#
41.00	23.20	53.98#
42.00	18.00	30.76#

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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	7.940	152	148676	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.745	136	707949	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	512628	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	1092578	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	828626	20.000	ng/ul	# 0.00
88) Perylene-d12	23.856	264	682495	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	14218m	4.170	ng/uL	0.00
4) Pyridine-d5	3.799	84	64834	6.760	ng/ul	0.01
7) Phenol-d5	7.098	99	80827	5.977	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	256997	31.671	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	243832	24.346	ng/ul	0.00
15) 4-Methylphenol-d8	8.645	113	172232	15.372	ng/ul	0.00
21) Nitrobenzene-d5	9.098	128	175270	31.128	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	174294	28.343	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	320059	26.735	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	360578	21.006	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	1352866	33.926	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1485721	30.785	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	23188m	3.159	ng/ul	0.00
60) Fluorene-d10	15.563	176	1186881	34.601	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.680	200	123324	17.851	ng/ul	0.00
73) Anthracene-d10	17.421	188	1901468	35.973	ng/ul	0.00
81) Pyrene-d10	19.651	212	2213339	46.968	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	1436977	40.161	ng/ul	0.00

Target Compounds Qvalue

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

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