

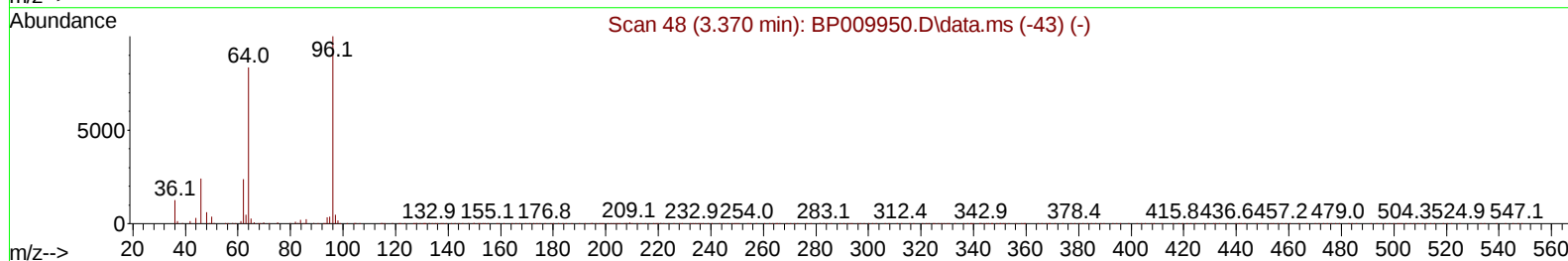
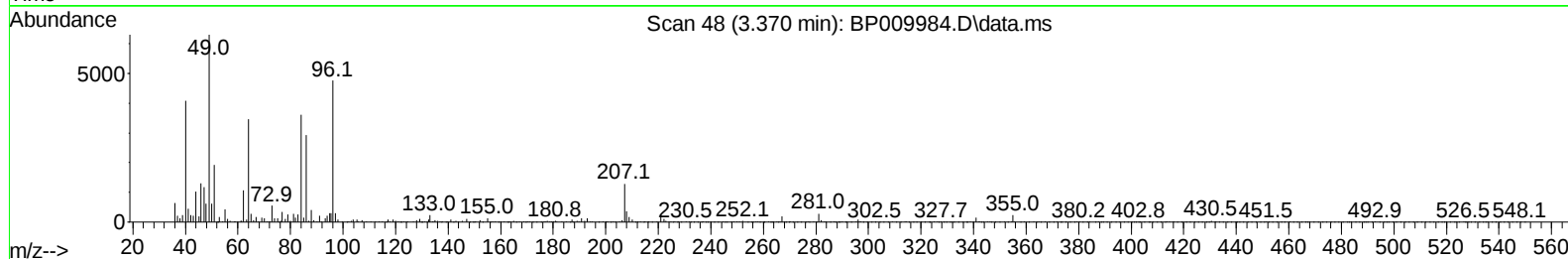
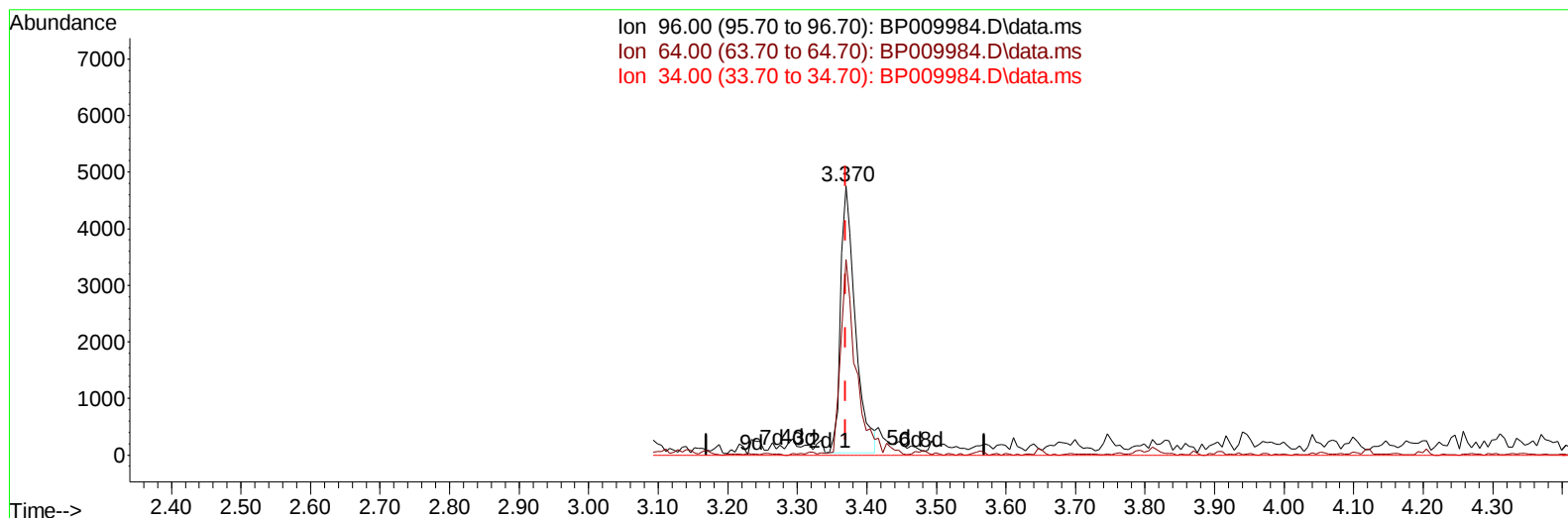
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009984.D
 Acq On : 20 Apr 2022 19:31
 Operator : CG/JU
 Sample : N2472-03
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J43

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/21/2022
 Supervised By :Yogesh Patel 04/25/2022

Quant Time: Apr 21 01:08:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 20 01:32:14 2022
 Response via : Initial Calibration



TIC: BP009984.D\data.ms

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

3.370min (-0.000) 2.39 ng/uL

response 6968

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	72.75#
34.00	0.00	0.00
0.00	0.00	0.00

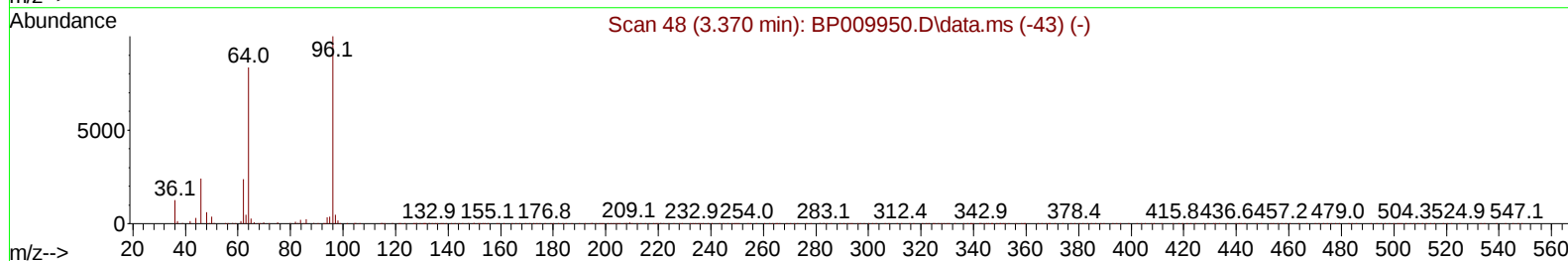
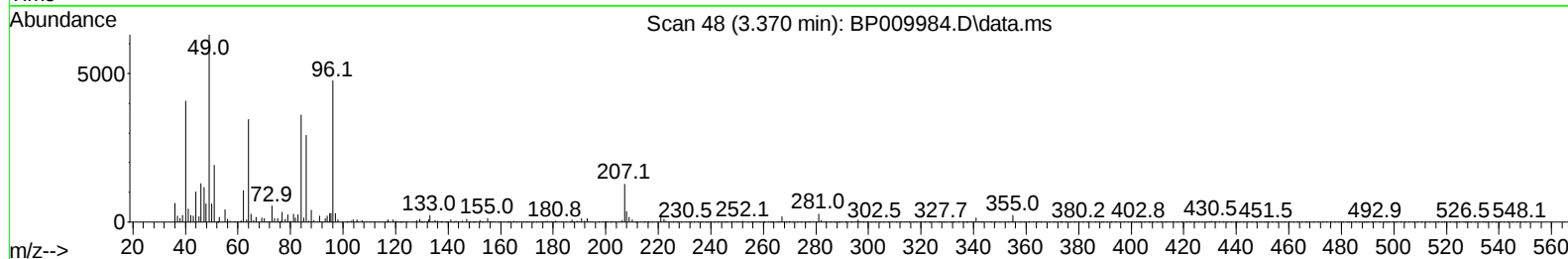
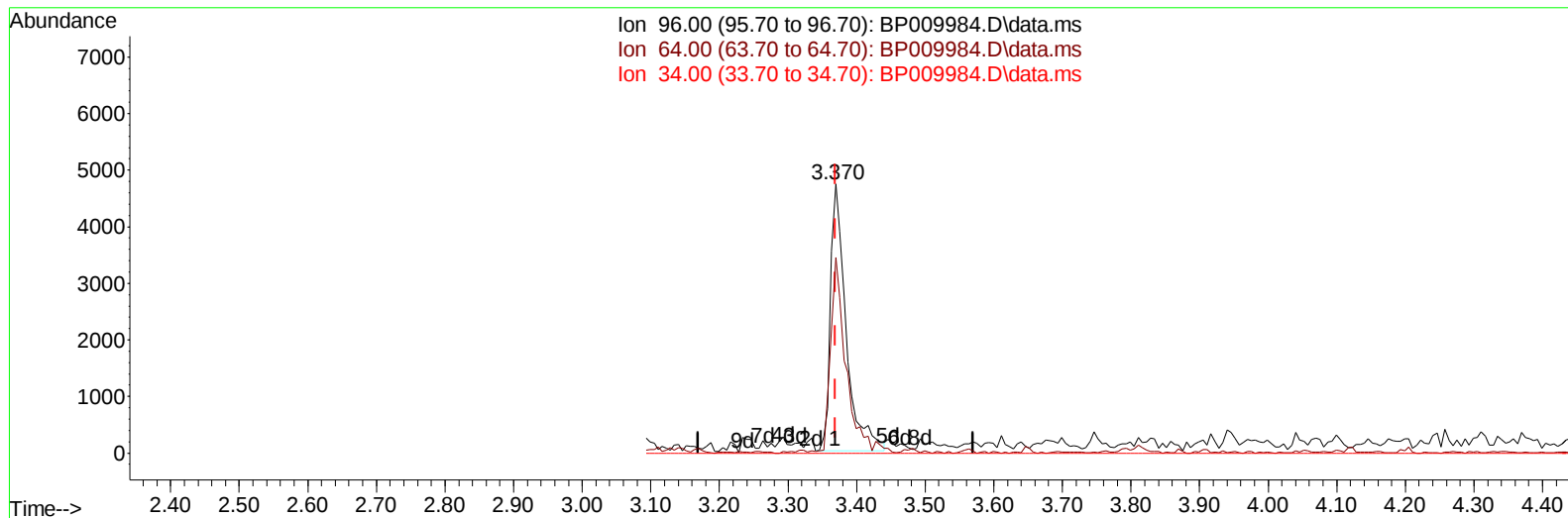
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TIC: BP009984.D\data.ms

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

3.370min (-0.000) 2.55 ng/uL m

response 7436

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	72.75#
34.00	0.00	0.00
0.00	0.00	0.00

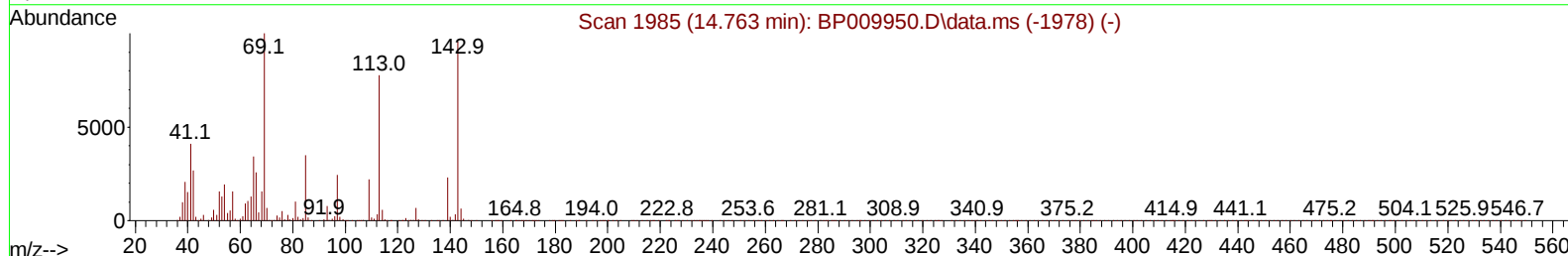
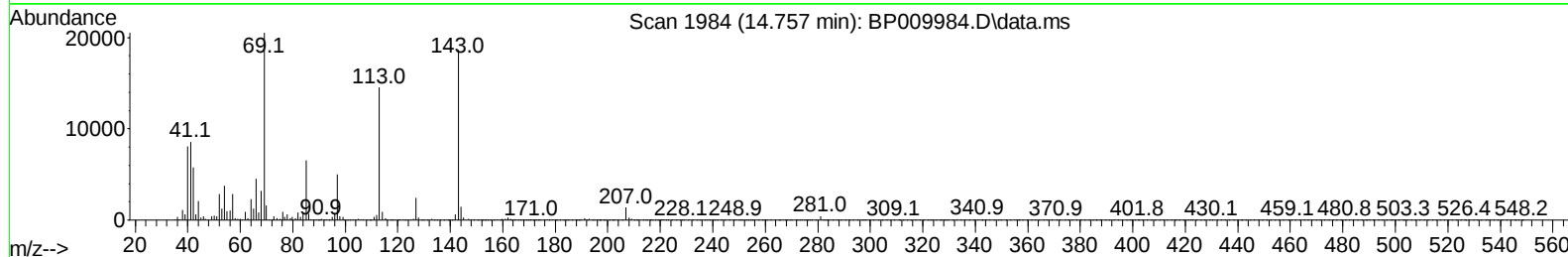
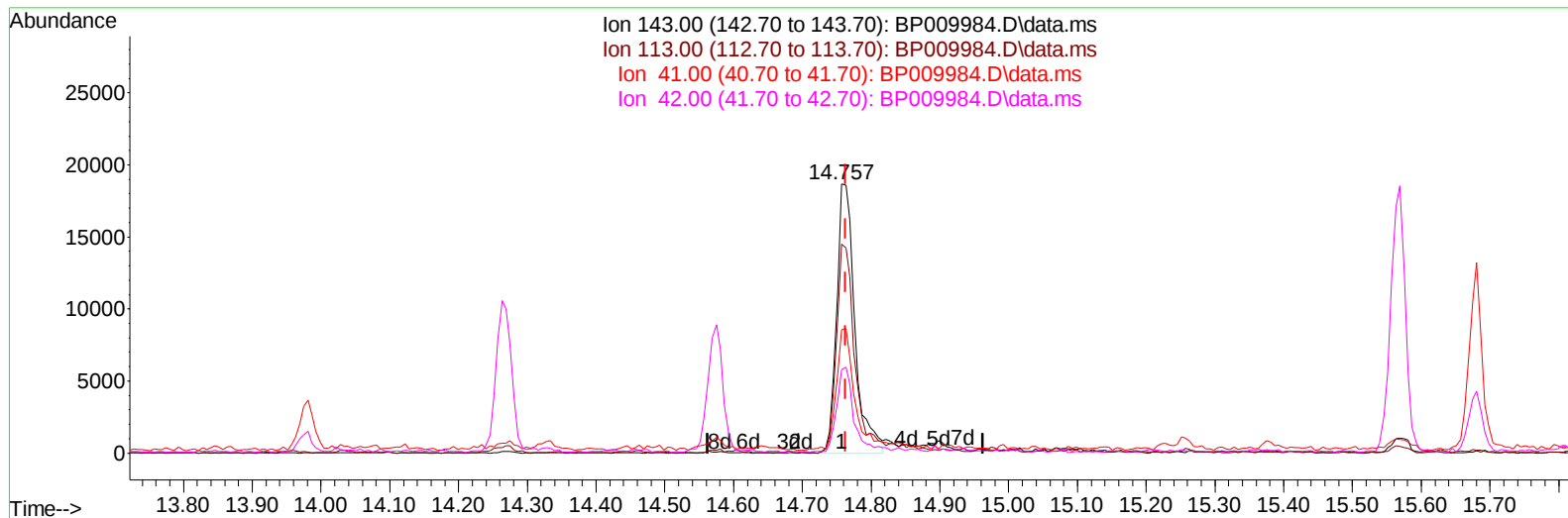
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TIC: BP009984.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (-0.006) 5.73 ng/u1

response 34378

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.79#
41.00	23.20	45.99#
42.00	18.00	31.09#

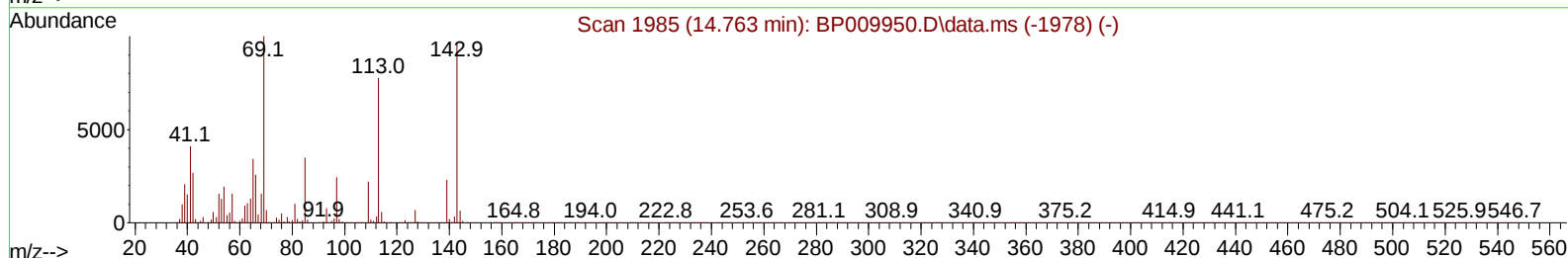
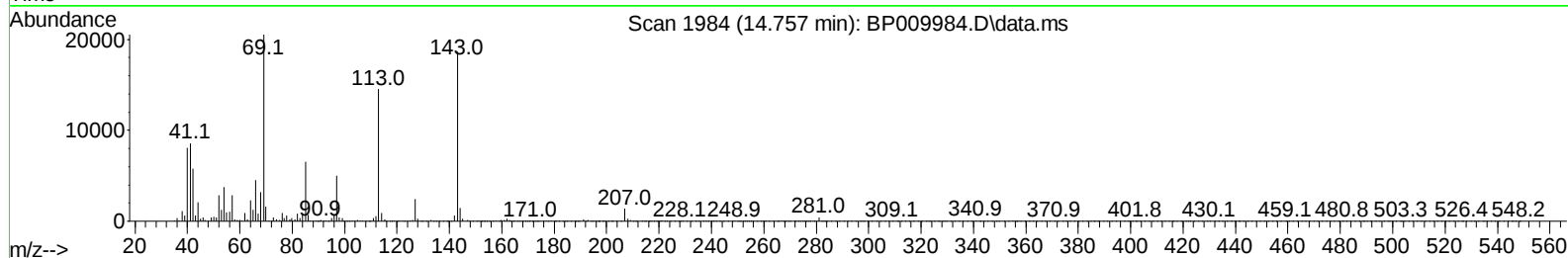
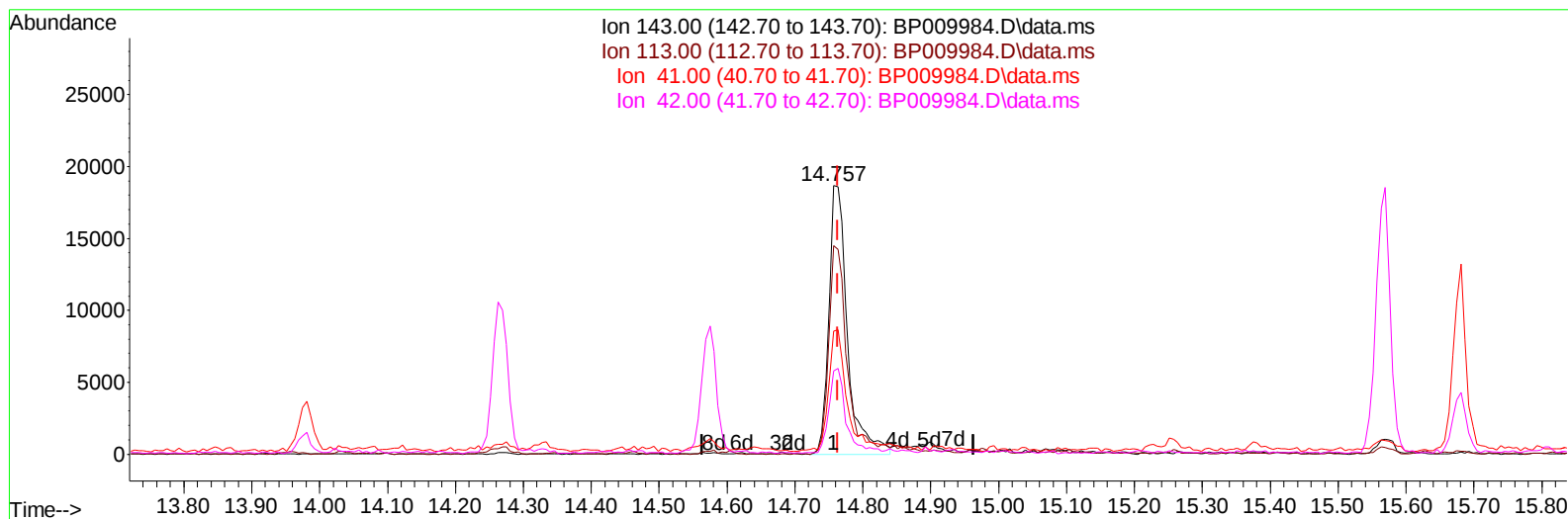
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 ALS Vial : 37 Sample Multiplier: 1

Instrument :
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TIC: BP009984.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (-0.006) 5.90 ng/ul m

response 35401

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.79#
41.00	23.20	45.99#
42.00	18.00	31.09#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	130037	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	590107	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	421176	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	916393	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	914858	20.000	ng/ul	# 0.00
88) Perylene-d12	23.863	264	810507	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	7436m	2.552	ng/uL	0.00
4) Pyridine-d5	3.793	84	44771	5.573	ng/ul	0.00
7) Phenol-d5	7.105	99	70404	6.167	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	245765	36.160	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	228191	26.498	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	152742	16.027	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	165011	38.777	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	177934	38.139	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	291682	29.877	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	357365	25.701	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1224526	37.013	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	1357673	34.639	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	35401m	5.900	ng/ul	0.00
60) Fluorene-d10	15.569	176	1056976	37.978	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	185878	34.240	ng/ul	0.00
73) Anthracene-d10	17.428	188	1810797	41.271	ng/ul	0.00
81) Pyrene-d10	19.657	212	2209092	44.213	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	1771780	41.977	ng/ul	0.00

Target Compounds Qvalue

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

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