

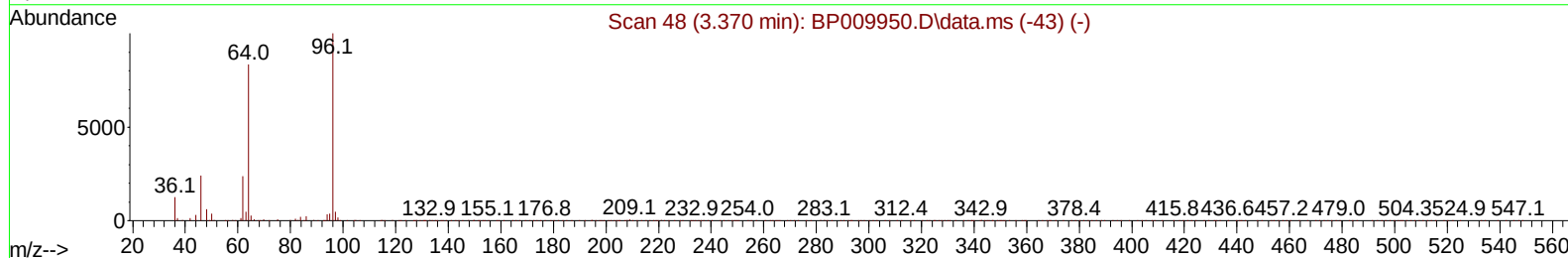
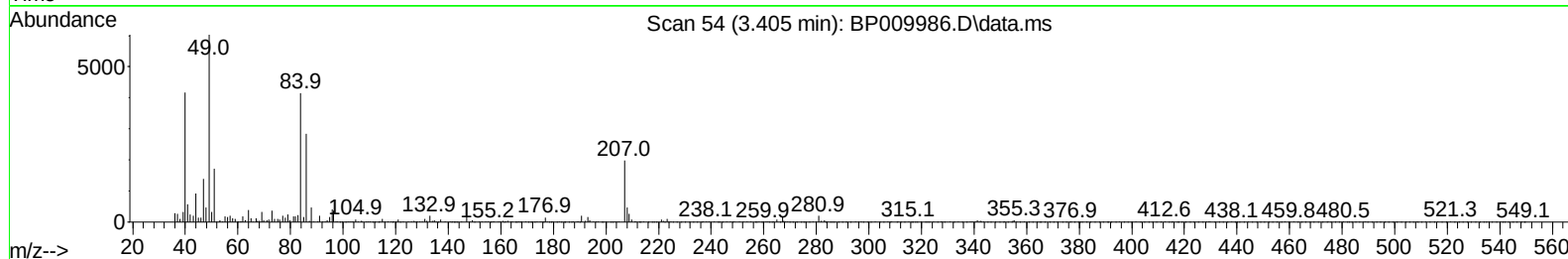
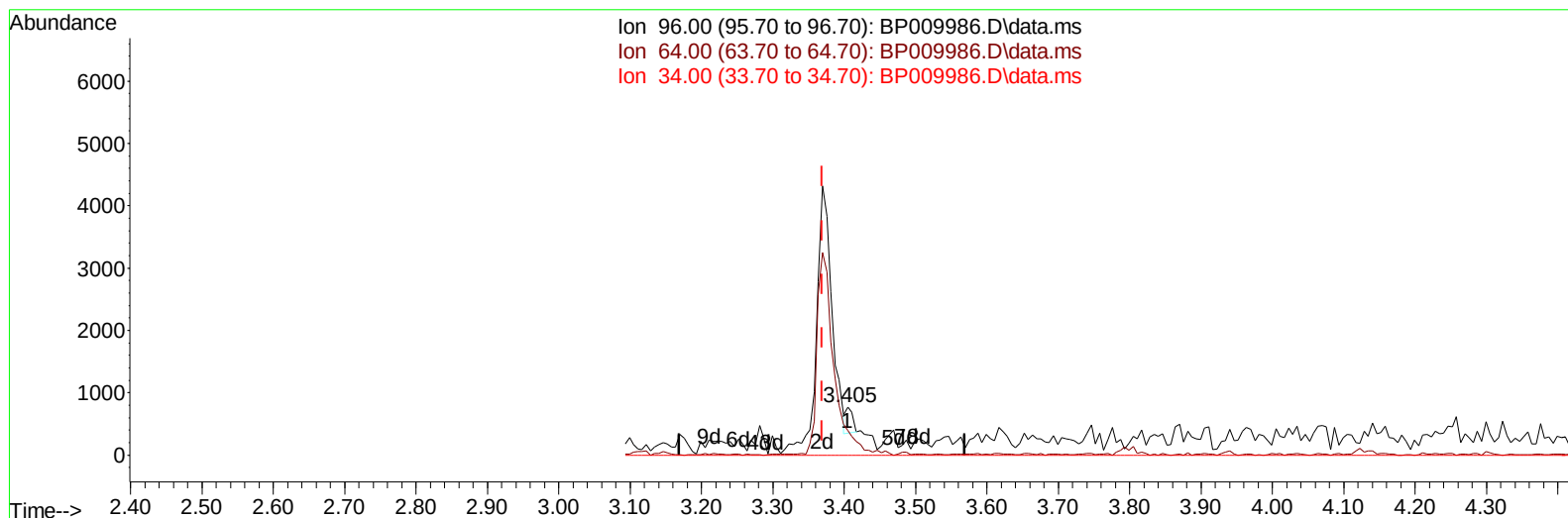
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009986.D
 Acq On : 20 Apr 2022 20:38
 Operator : CG/JU
 Sample : N2472-12
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J32

Manual IntegrationsAPPROVED

Quant Time: Apr 21 01:08:54 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

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



TIC: BP009986.D\data.ms

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

3.405min (+ 0.035) 0.12 ng/uL

response 269

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	50.32
34.00	0.00	0.00
0.00	0.00	0.00

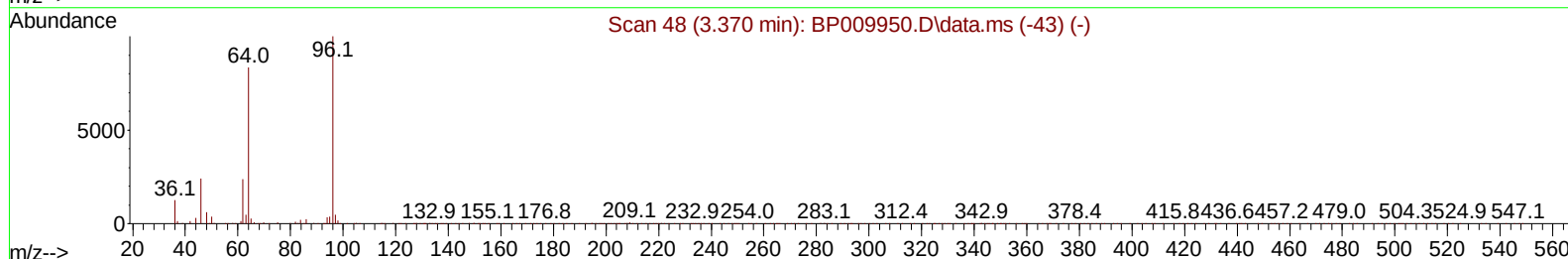
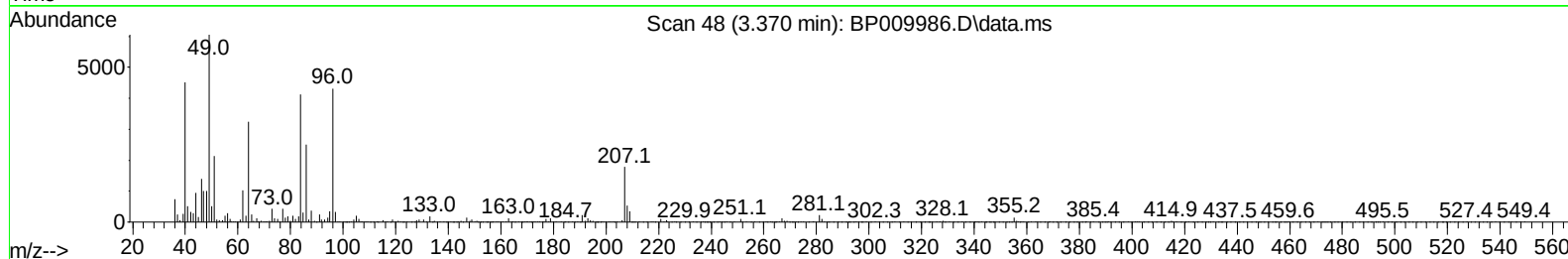
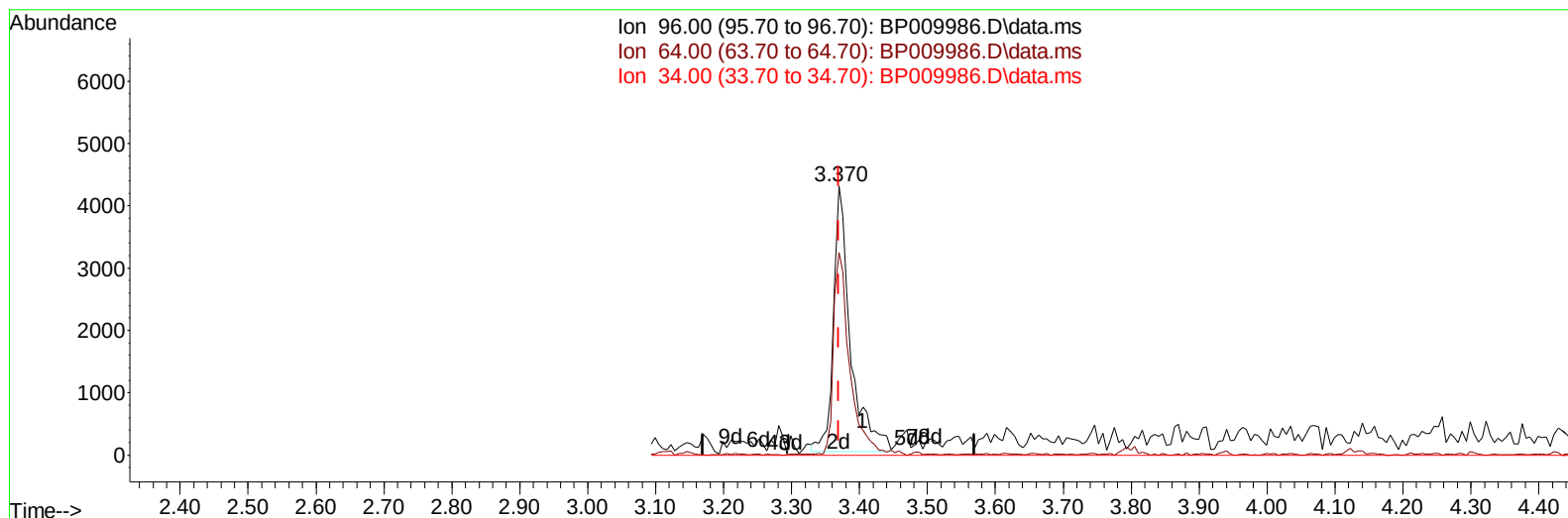
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(3) 1,4-Dioxane-d8 (S)

3.370min (0.000) 3.28 ng/uL m

response 7457

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	75.29#
34.00	0.00	0.00
0.00	0.00	0.00

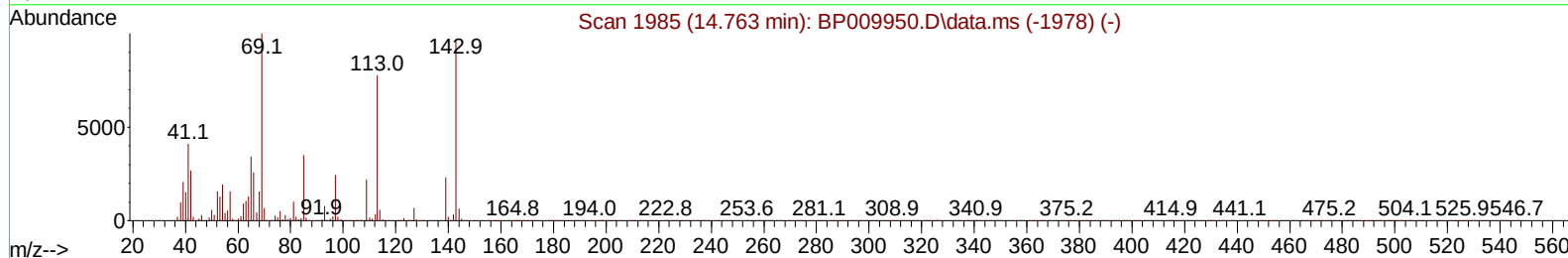
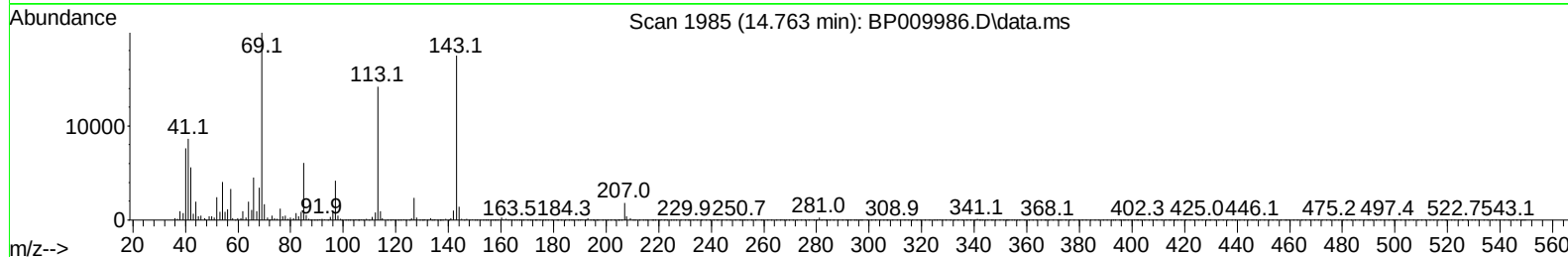
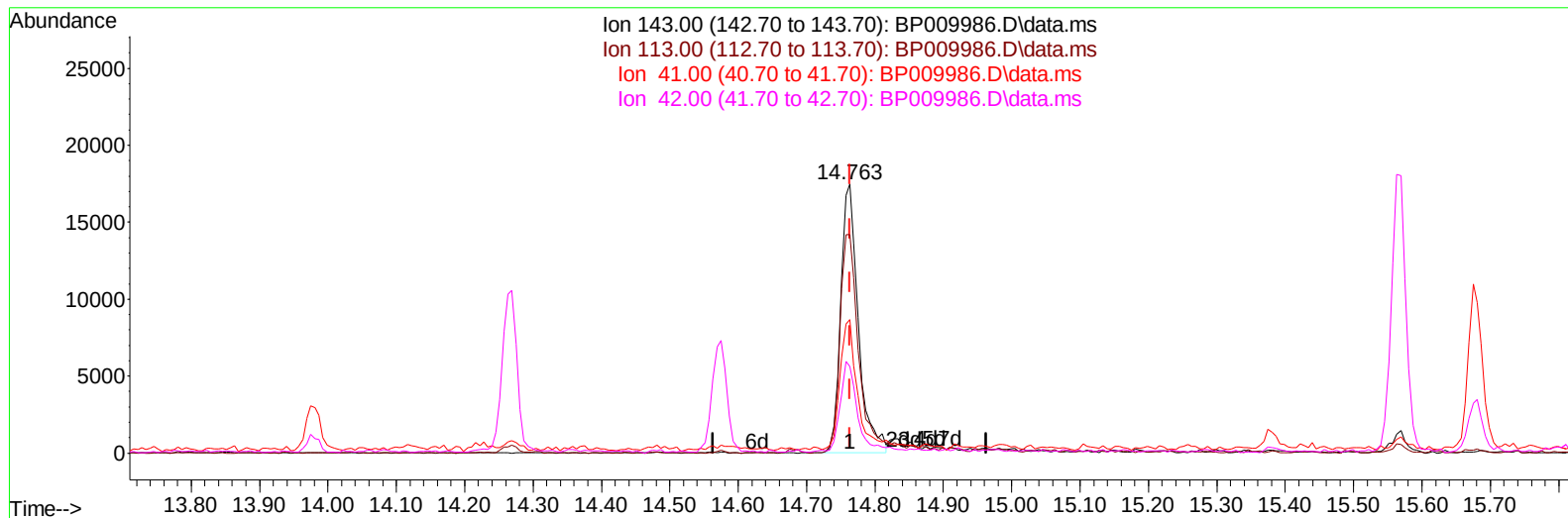
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TIC: BP009986.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (0.000) 6.56 ng/u1

response 31269

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	81.28#
41.00	23.20	49.55#
42.00	18.00	32.21#

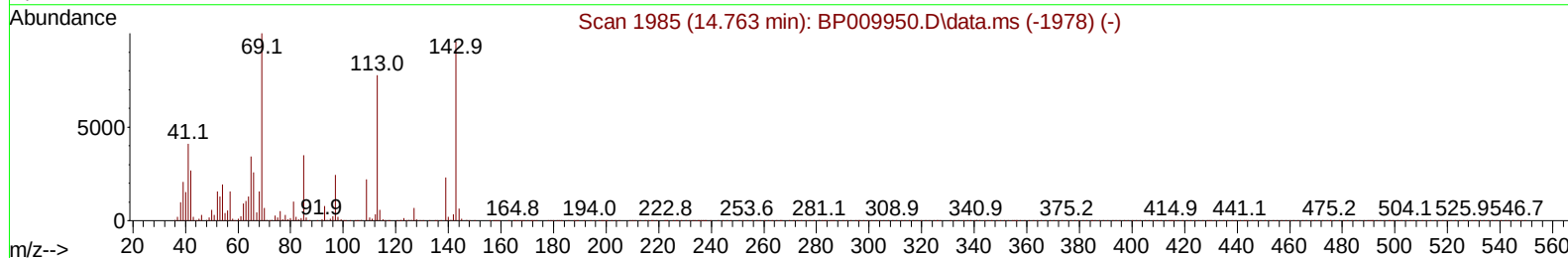
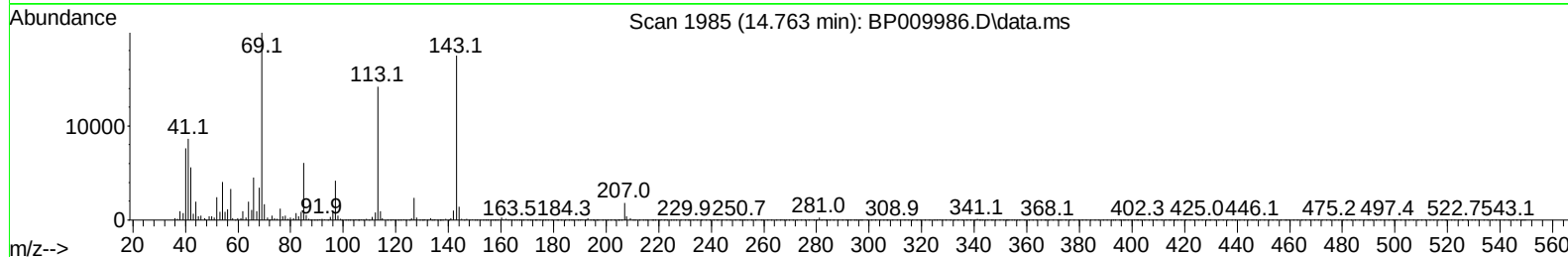
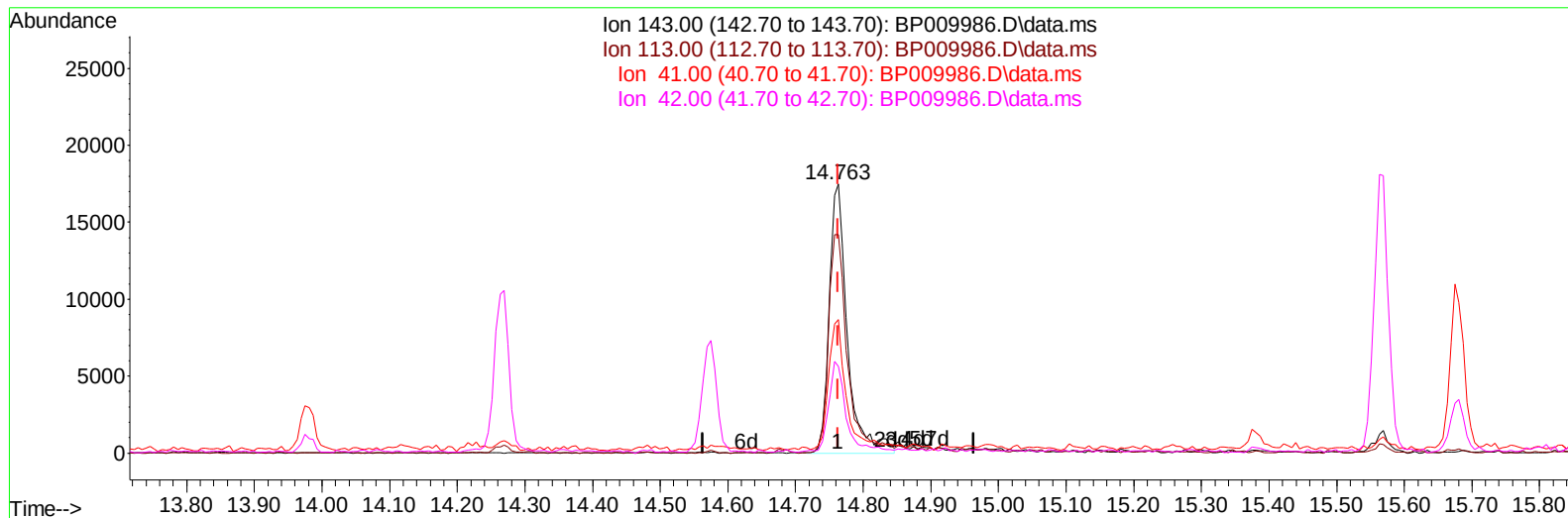
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Instrument :
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TIC: BP009986.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (0.000) 6.85 ng/ul m

response 32683

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	81.28#
41.00	23.20	49.55#
42.00	18.00	32.21#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	101396	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	466550	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.575	164	334713	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	738654	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	759014	20.000	ng/ul	# 0.00
88) Perylene-d12	23.857	264	745511	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	7457m	3.282	ng/uL	0.00
4) Pyridine-d5	3.793	84	54613	8.718	ng/ul	0.00
7) Phenol-d5	7.105	99	69498	7.807	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	236490	44.624	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	218979	32.611	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	146193	19.673	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	156760	46.593	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	167261	45.345	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	271436	35.166	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	362695	32.992	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1174101	44.656	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	1301316	41.777	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	32683m	6.854	ng/ul	0.00
60) Fluorene-d10	15.569	176	1024953	46.340	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	172804	39.491	ng/ul	0.00
73) Anthracene-d10	17.428	188	1730293	48.925	ng/ul	0.00
81) Pyrene-d10	19.651	212	2096180	50.567	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	1869294	48.148	ng/ul	0.00

Target Compounds Qvalue

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

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