

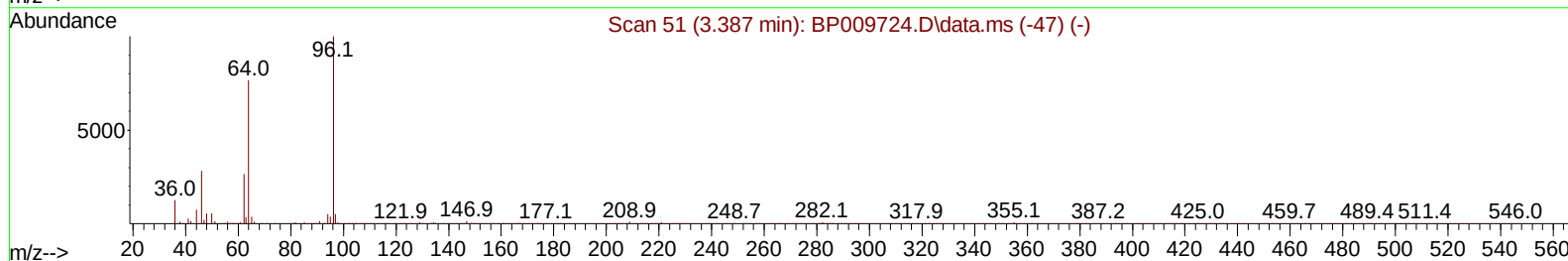
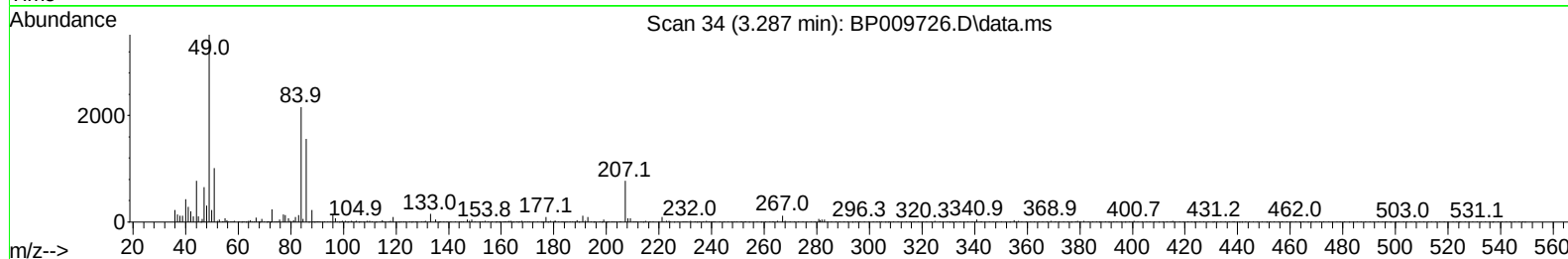
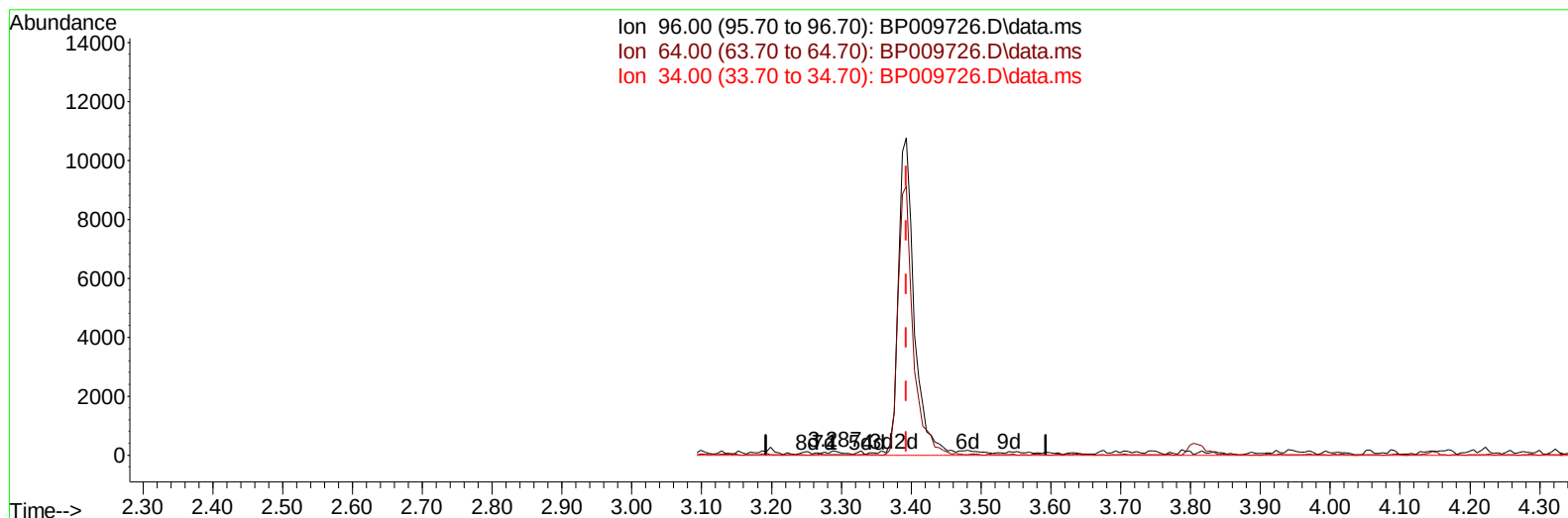
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
 Data File : BP009726.D
 Acq On : 08 Apr 2022 20:28
 Operator : CG/JU
 Sample : PB143934BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK934

Manual IntegrationsAPPROVED

Quant Time: Apr 08 22:30:21 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 19:16:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/11/2022
 Supervised By :Yogesh Patel 04/12/2022



TIC: BP009726.D\data.ms

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

3.287min (-0.106) 0.03 ng/uL

response 86

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	45.26
34.00	0.00	0.00
0.00	0.00	0.00

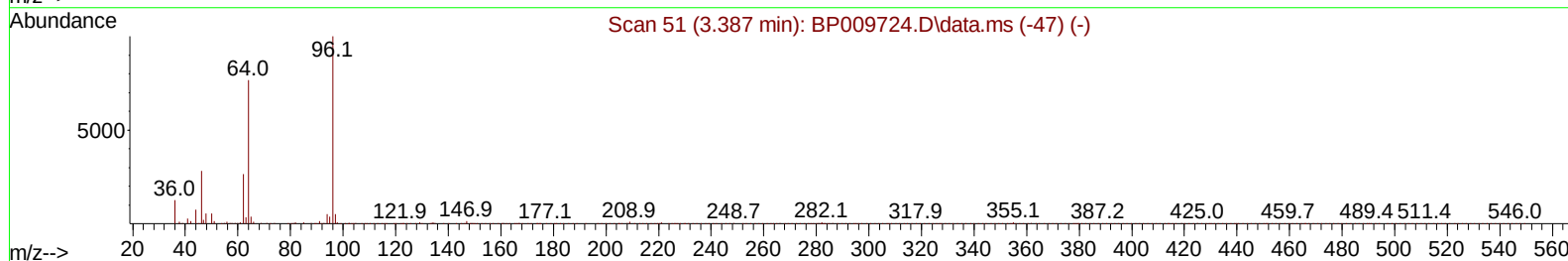
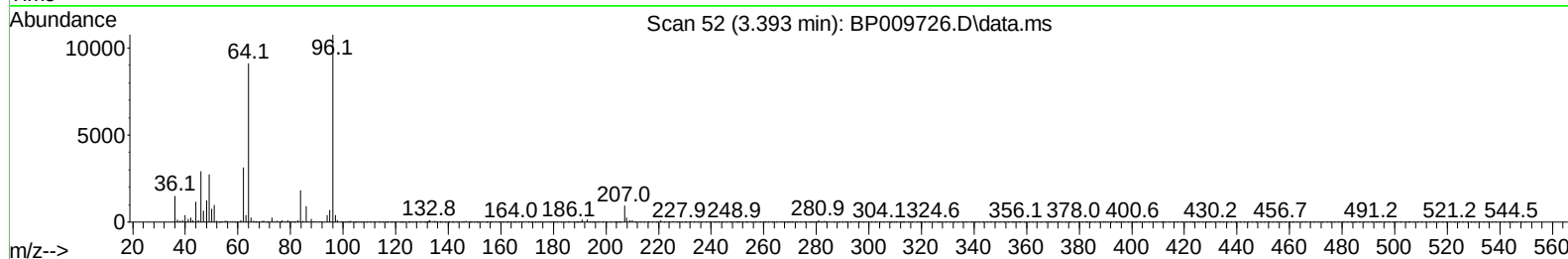
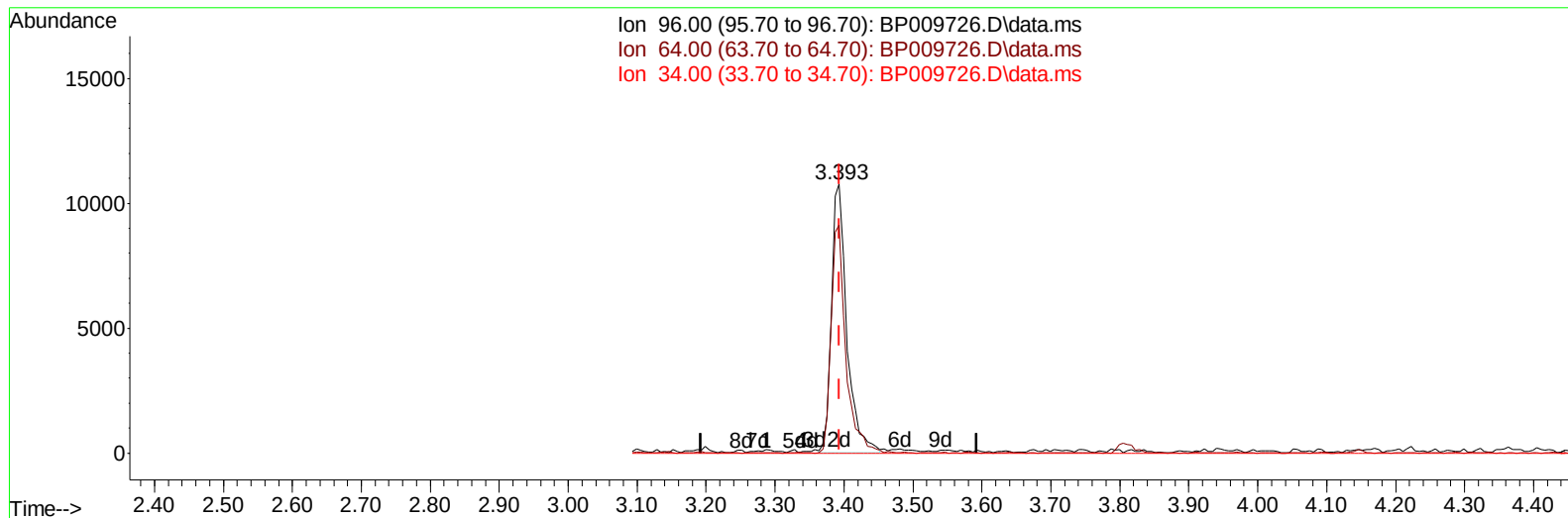
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TIC: BP009726.D\data.ms

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

3.393min (+ 0.000) 6.31 ng/uL m

response 16699

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	84.75#
34.00	0.00	0.00
0.00	0.00	0.00

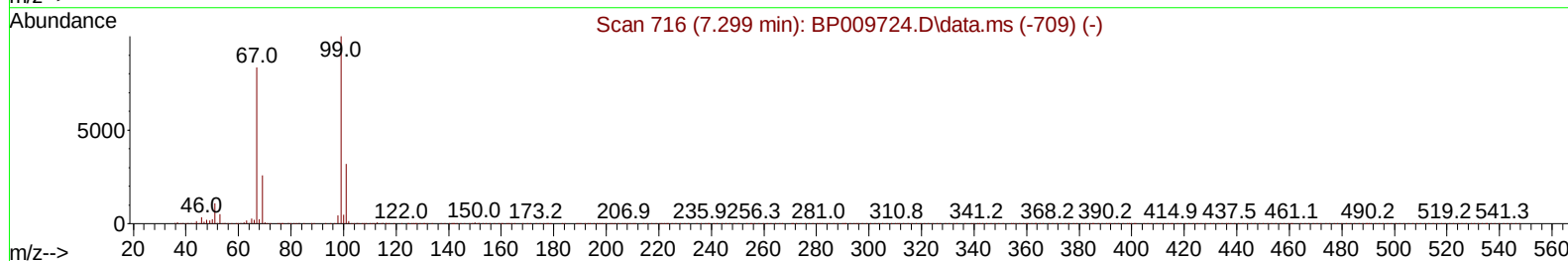
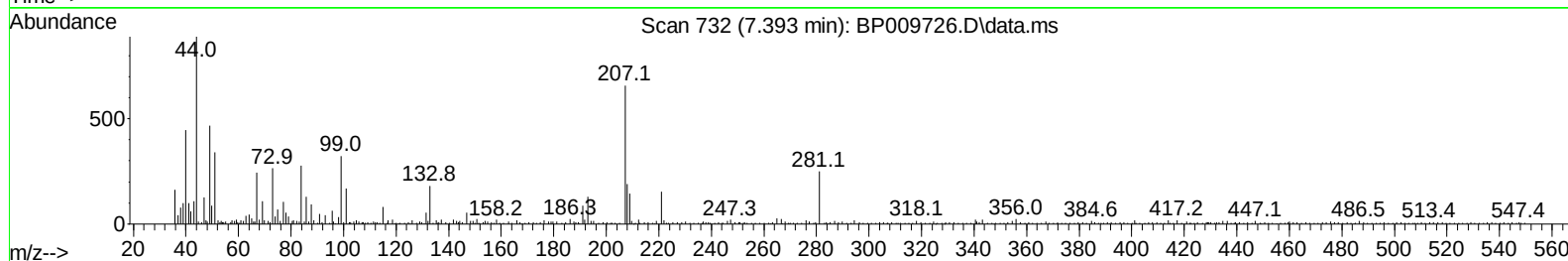
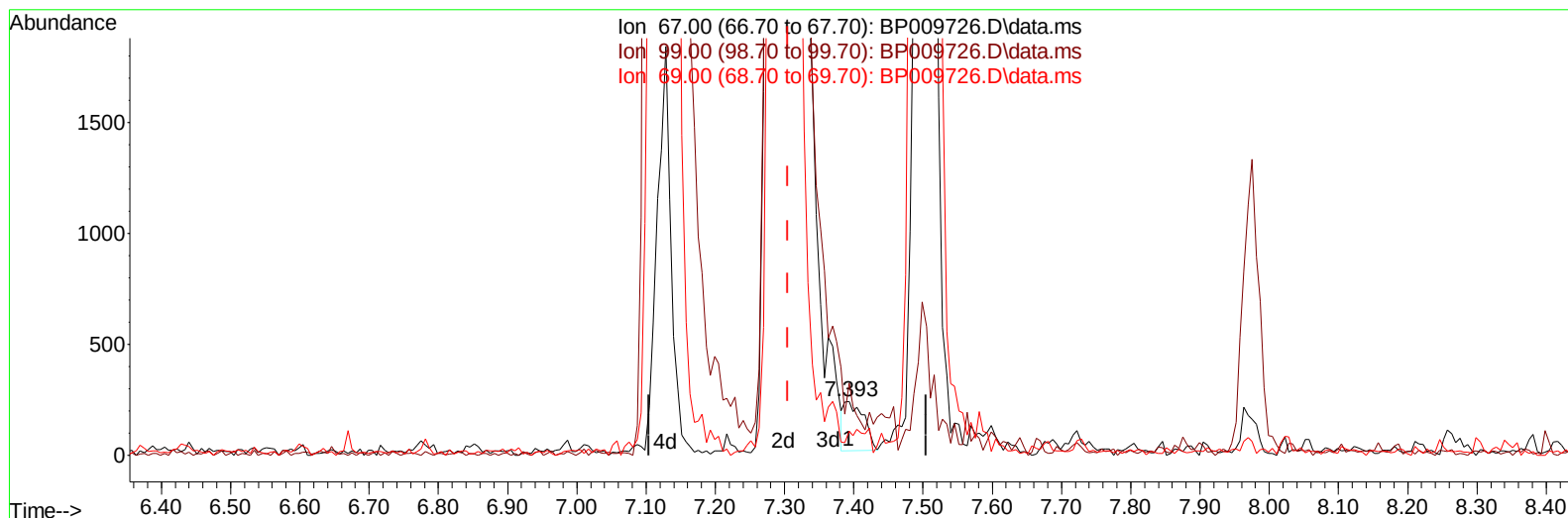
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TIC: BP009726.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.393min (+ 0.088) 0.07 ng/u1

response 447

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	133.33
69.00	40.10	44.86
0.00	0.00	0.00

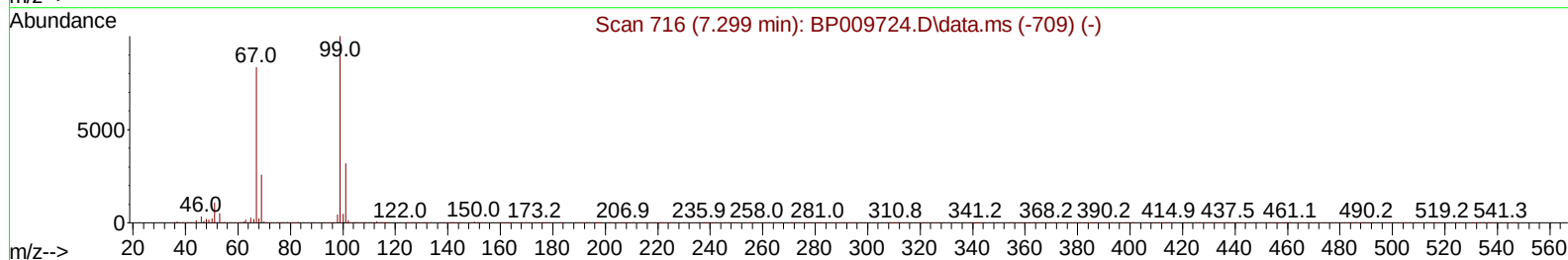
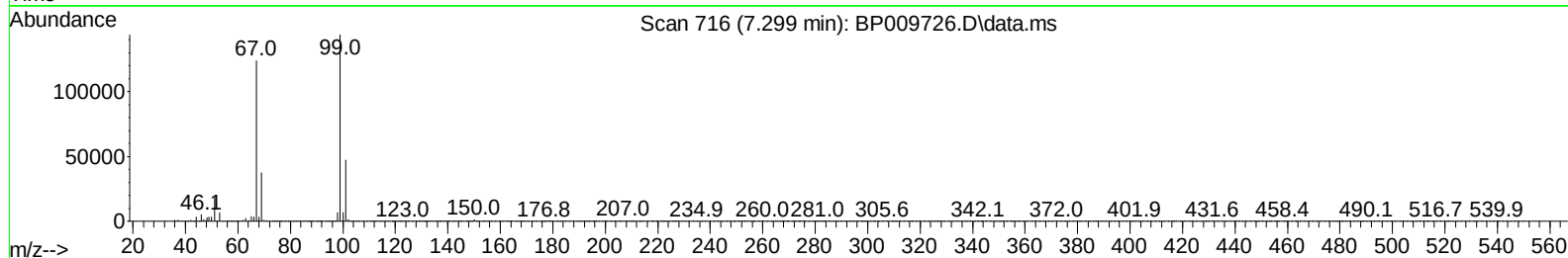
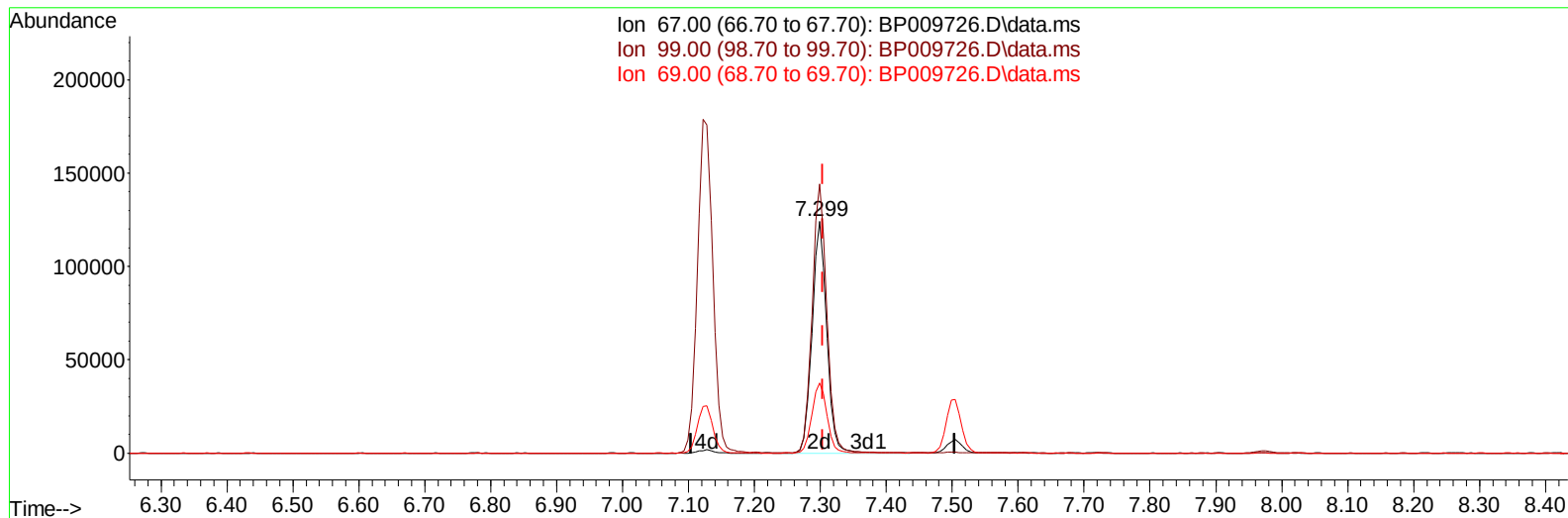
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TIC: BP009726.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.299min (-0.006) 31.06 ng/ul m

response 191898

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	116.02#
69.00	40.10	30.16#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	118194	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	517896	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.604	164	379088	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.357	188	846251	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.439	240	905475	20.000	ng/ul	# 0.00
88) Perylene-d12	23.916	264	884682	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	16699m	6.305	ng/uL	0.00
4) Pyridine-d5	3.805	84	227332	31.134	ng/ul	0.00
7) Phenol-d5	7.122	99	287548	27.711	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	191898m	31.064	ng/ul	0.00
11) 2-Chlorophenol-d4	7.505	132	231768	29.610	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	248116	28.643	ng/ul	0.00
21) Nitrobenzene-d5	9.134	128	122884	32.903	ng/ul	0.00
24) 2-Nitrophenol-d4	9.857	143	136167	33.256	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	244788	28.570	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	366548	30.037	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	949119	31.873	ng/ul	0.00
49) Acenaphthylene-d8	14.298	160	1037883	29.420	ng/ul	0.00
54) 4-Nitrophenol-d4	14.781	143	156452	28.971	ng/ul	0.00
60) Fluorene-d10	15.598	176	789853	31.531	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	142508	28.426	ng/ul	0.00
73) Anthracene-d10	17.457	188	1295614	31.976	ng/ul	0.00
81) Pyrene-d10	19.680	212	1599188	32.338	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.757	264	1537759	33.378	ng/ul	0.00

Target Compounds Qvalue

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

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