

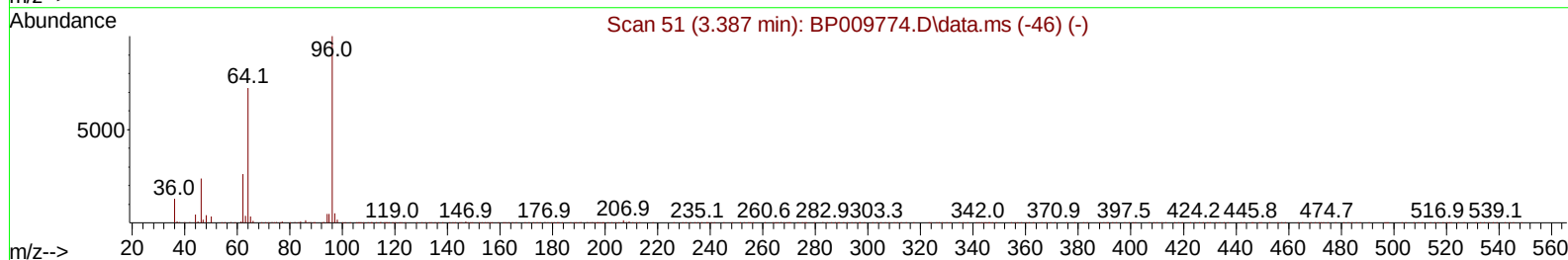
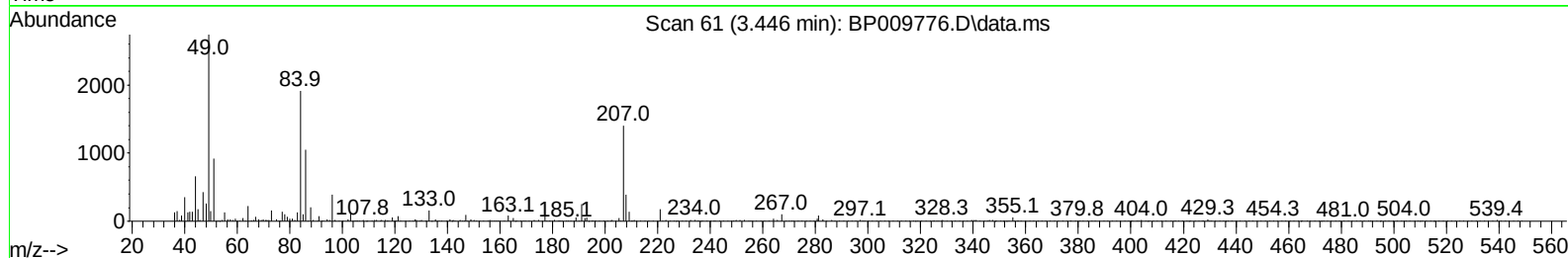
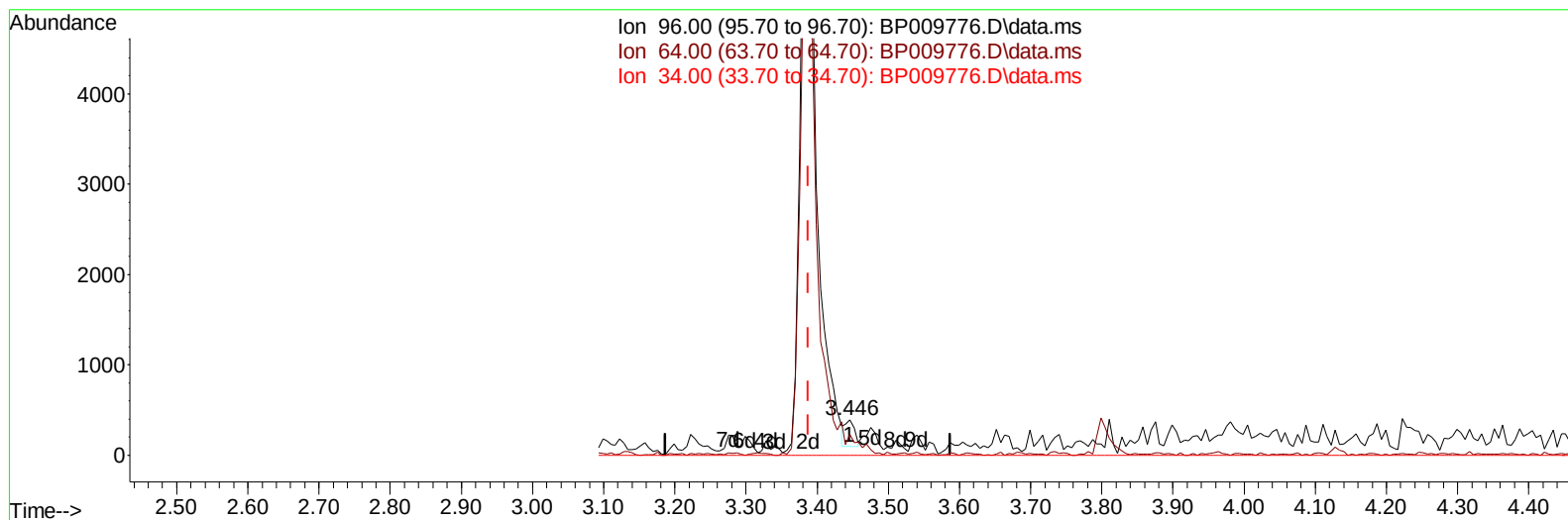
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009776.D
 Acq On : 12 Apr 2022 09:51
 Operator : CG/JU
 Sample : PB144054BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK054

Manual IntegrationsAPPROVED

Quant Time: Apr 13 02:34:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 13 02:15:22 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/13/2022
 Supervised By :Yogesh Patel 04/13/2022



TIC: BP009776.D\data.ms

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

3.446min (+ 0.059) 0.16 ng/uL

response 278

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	57.51
34.00	0.00	0.00
0.00	0.00	0.00

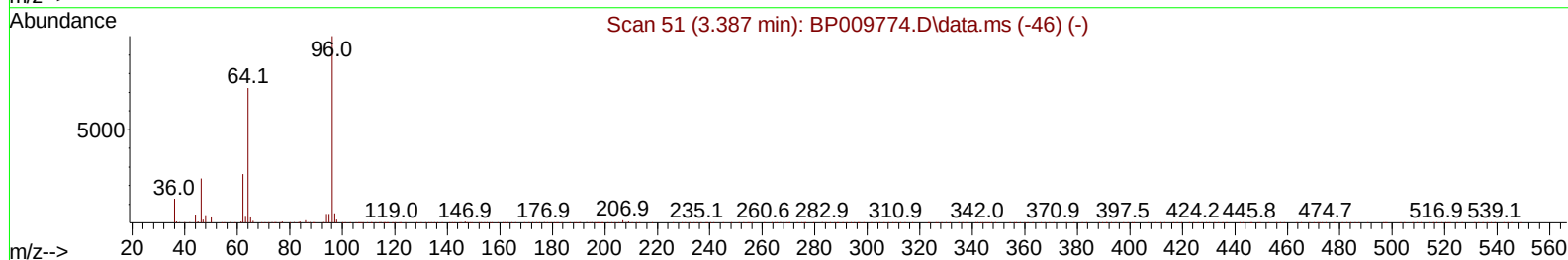
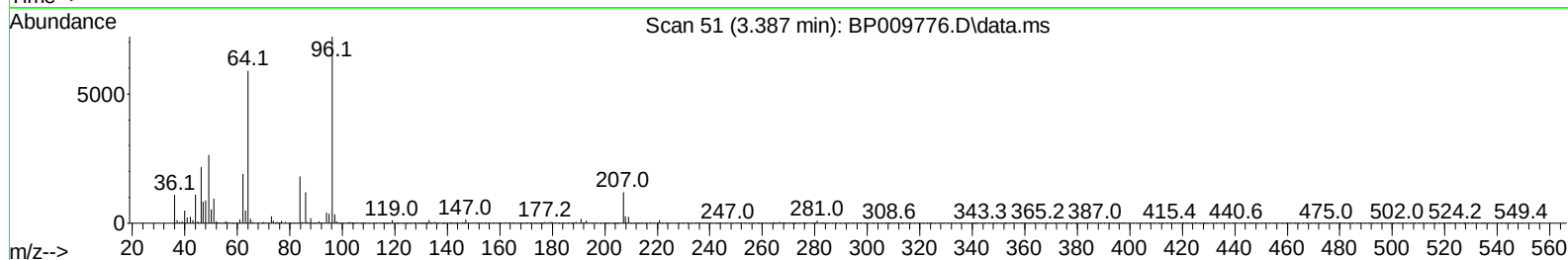
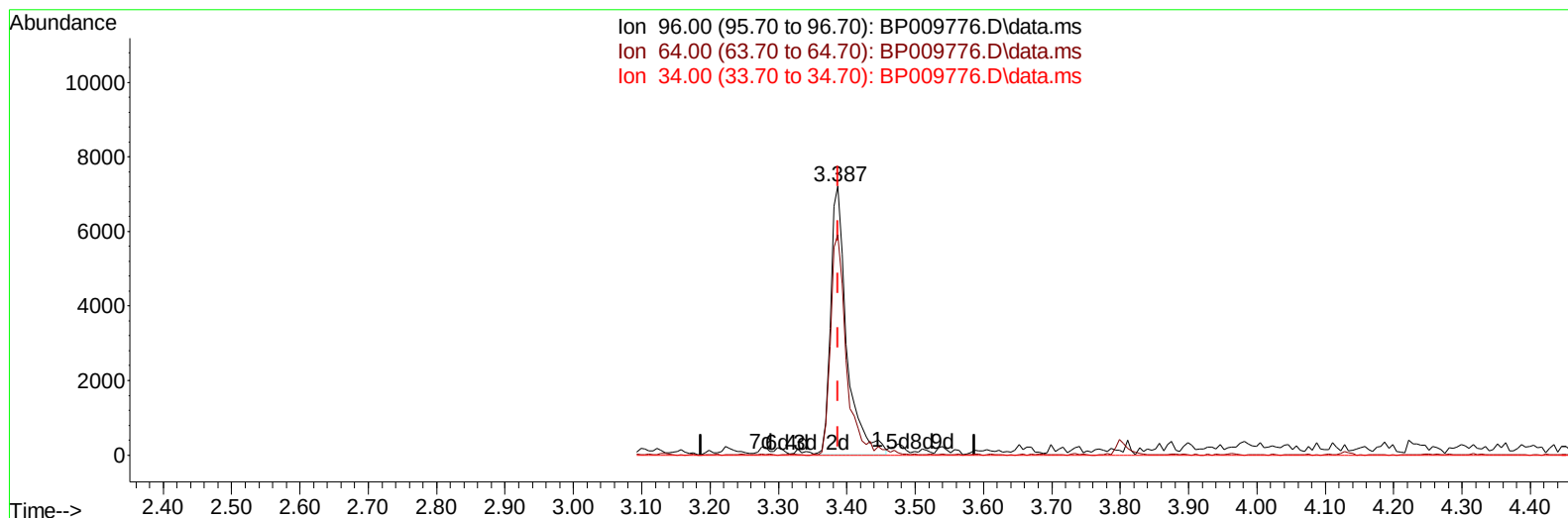
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(3) 1,4-Dioxane-d8 (S)

3.387min (-0.000) 6.64 ng/uL m

response 11796

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	81.85#
34.00	0.00	0.00
0.00	0.00	0.00

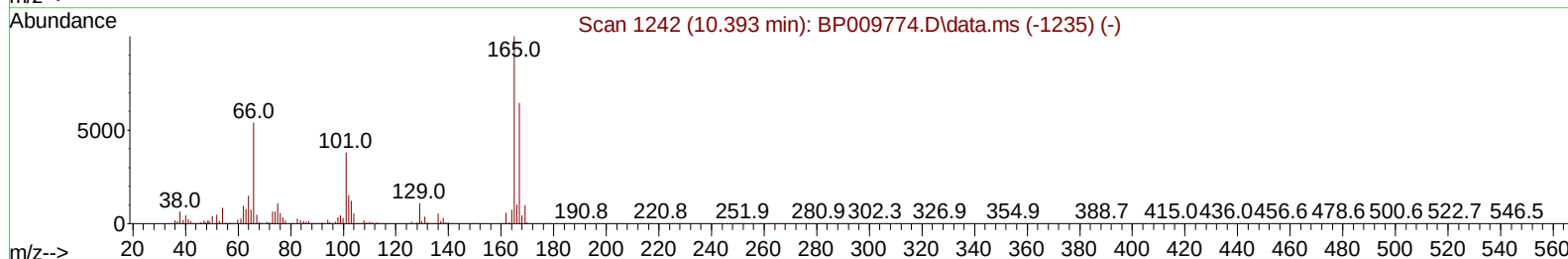
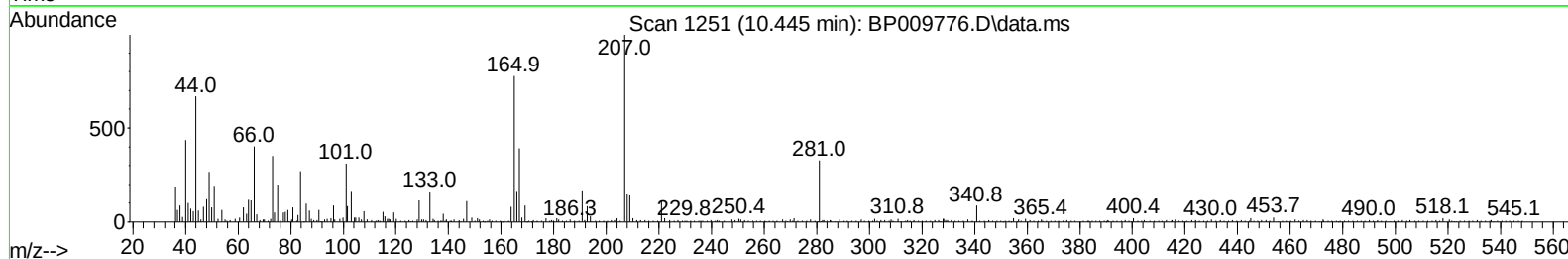
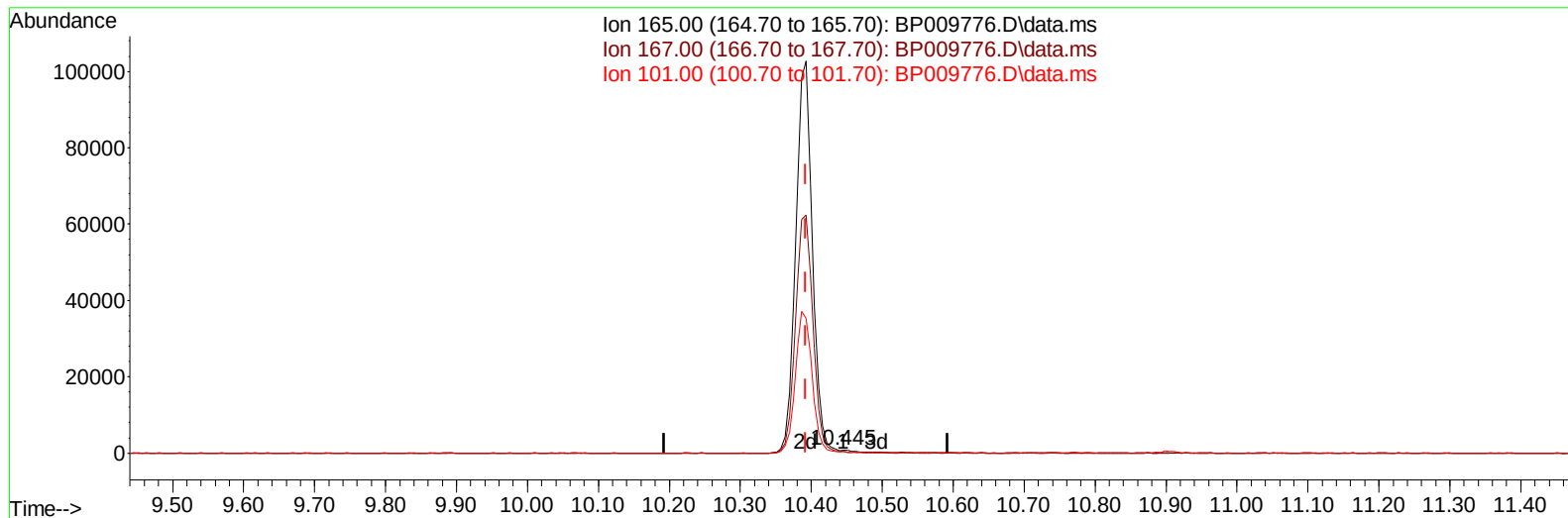
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TIC: BP009776.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.445min (+ 0.053) 0.14 ng/ul

response 822

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	50.32
101.00	49.80	50.97
0.00	0.00	0.00

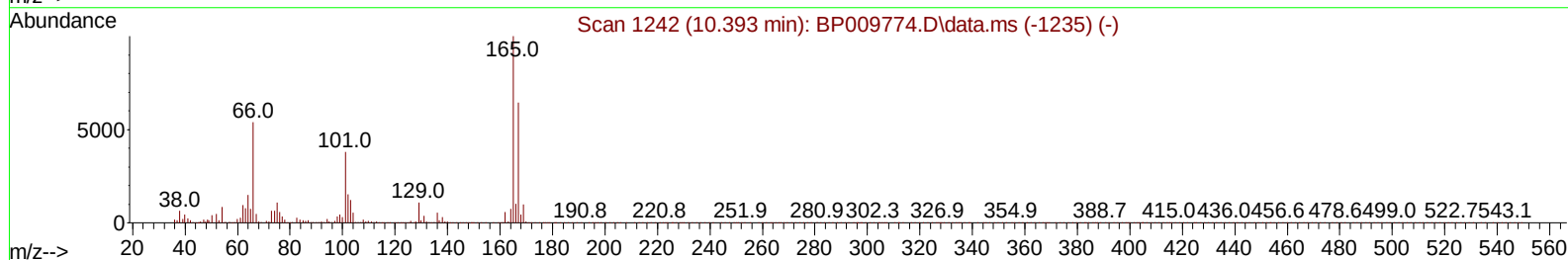
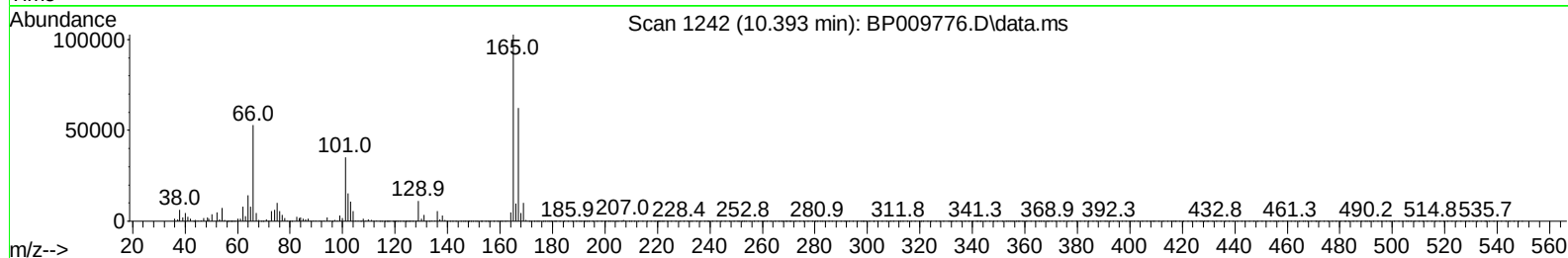
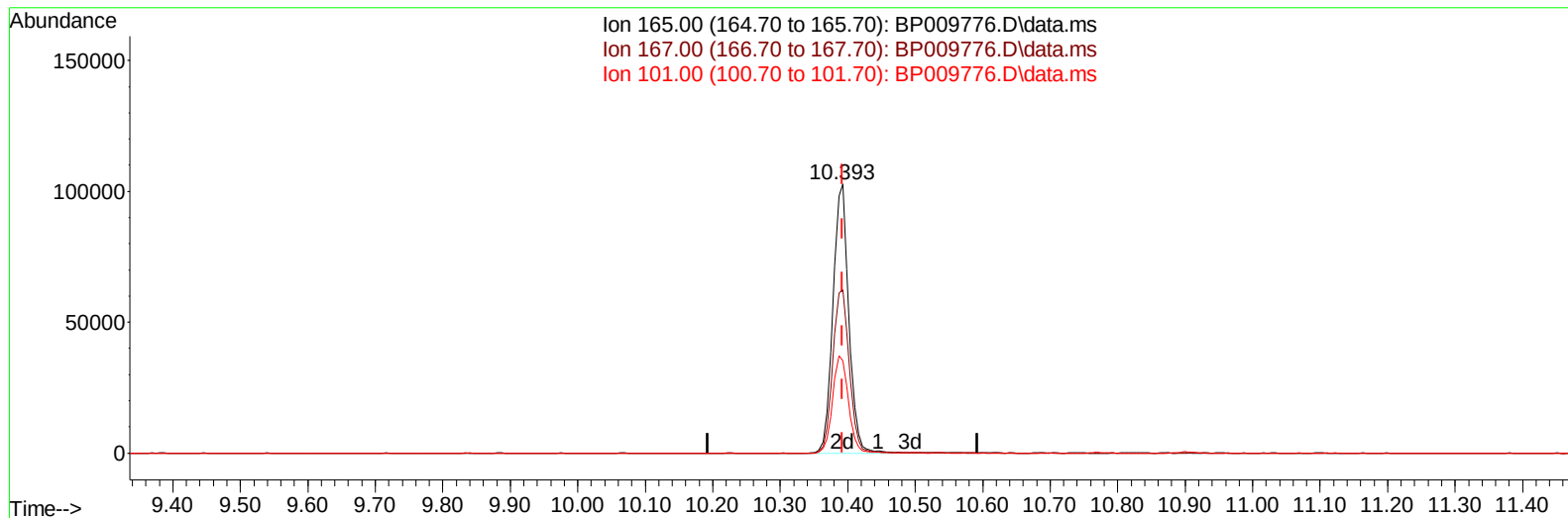
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TIC: BP009776.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.393min (-0.000) 29.11 ng/ul m

response 167353

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	60.70
101.00	49.80	34.33#
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	7.969	152	79252	20.000	ng/ul	0.00
20) Naphthalene-d8	10.775	136	347539	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.598	164	255199	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.351	188	568419	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.427	240	636267	20.000	ng/ul	# 0.00
88) Perylene-d12	23.898	264	662485	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	11796m	6.643	ng/uL	0.00
4) Pyridine-d5	3.805	84	158376	32.348	ng/ul	0.00
7) Phenol-d5	7.116	99	199237	28.635	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	136008	32.834	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	160732	30.625	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	172794	29.749	ng/ul	0.00
21) Nitrobenzene-d5	9.122	128	86390	34.470	ng/ul	0.00
24) 2-Nitrophenol-d4	9.851	143	91720	33.381	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	167353m	29.106	ng/ul	0.00
31) 4-Chloroaniline-d4	10.904	131	260634	31.827	ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	663285	33.088	ng/ul	0.00
49) Acenaphthylene-d8	14.292	160	706529	29.750	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	110085	30.281	ng/ul	0.00
60) Fluorene-d10	15.592	176	552509	32.763	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	92626	27.507	ng/ul	0.00
73) Anthracene-d10	17.451	188	928285	34.109	ng/ul	0.00
81) Pyrene-d10	19.674	212	1160274	33.389	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.739	264	1191249	34.529	ng/ul	0.00

Target Compounds Qvalue

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

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