

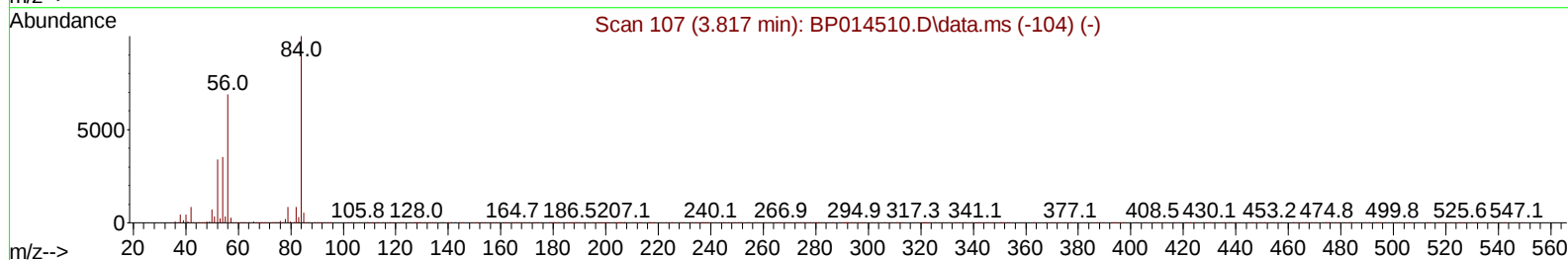
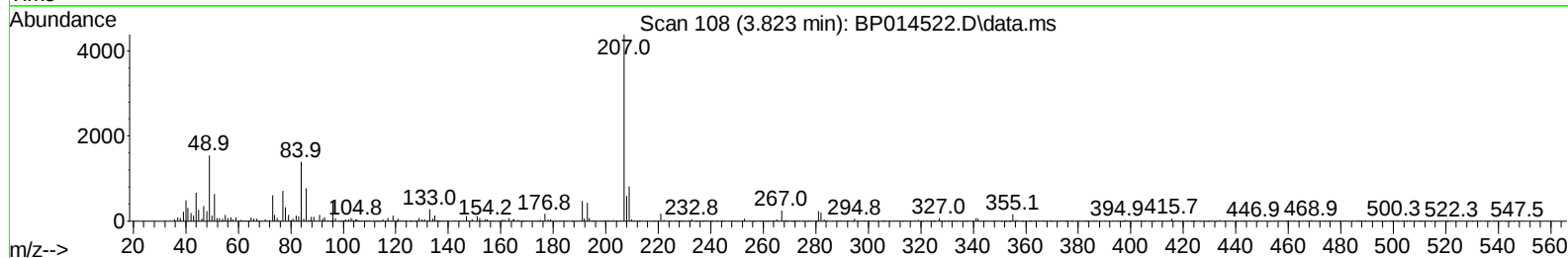
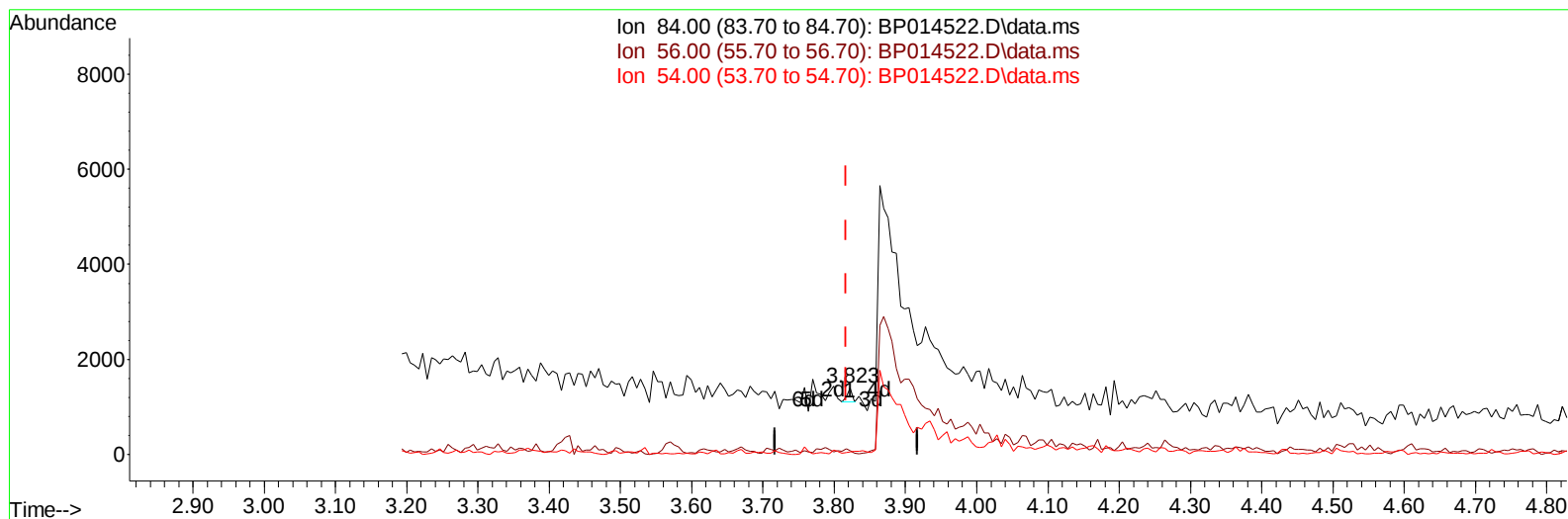
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
 Data File : BP014522.D
 Acq On : 04 Apr 2023 11:45
 Operator : CG/JU
 Sample : 02092-19DL 10X
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB991DL

Manual IntegrationsAPPROVED

Quant Time: Apr 04 23:09:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 04 01:34:36 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/05/2023
 Supervised By :Jagrut Upadhyay 04/05/2023



TIC: BP014522.D\data.ms

(4) Pyridine-d5 (S)

3.823min (+ 0.006) 0.01 ng/u1

response 142

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	6.71#
54.00	32.60	3.79#
0.00	0.00	0.00

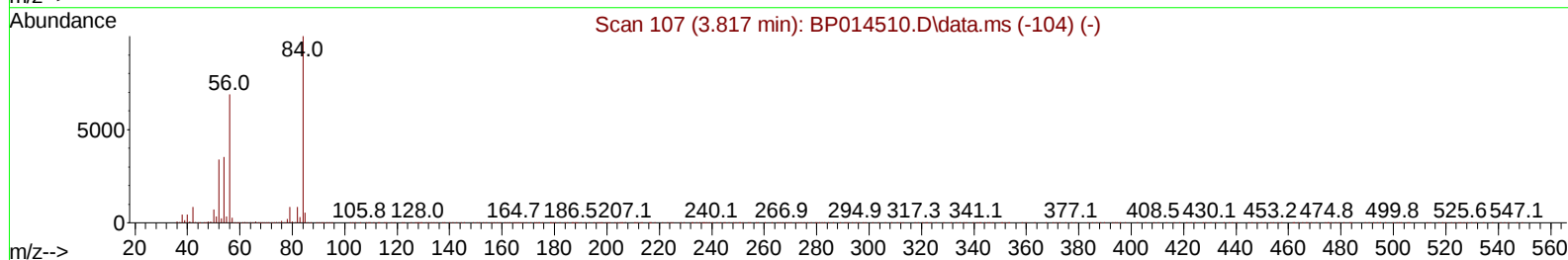
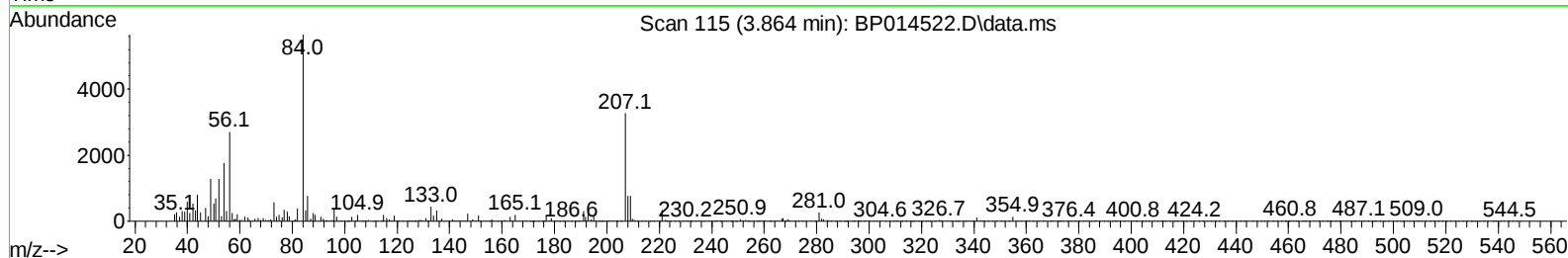
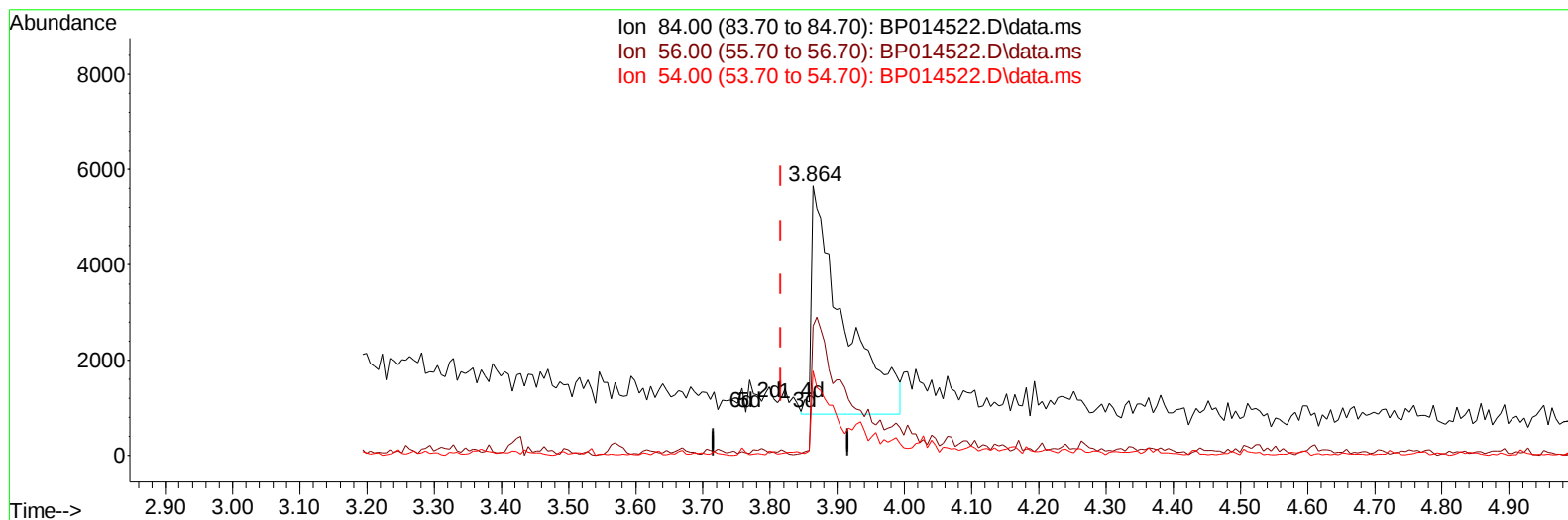
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(4) Pyridine-d5 (S)

3.864min (+ 0.047) 1.67 ng/ul m

response 15879

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	47.90#
54.00	32.60	31.16
0.00	0.00	0.00

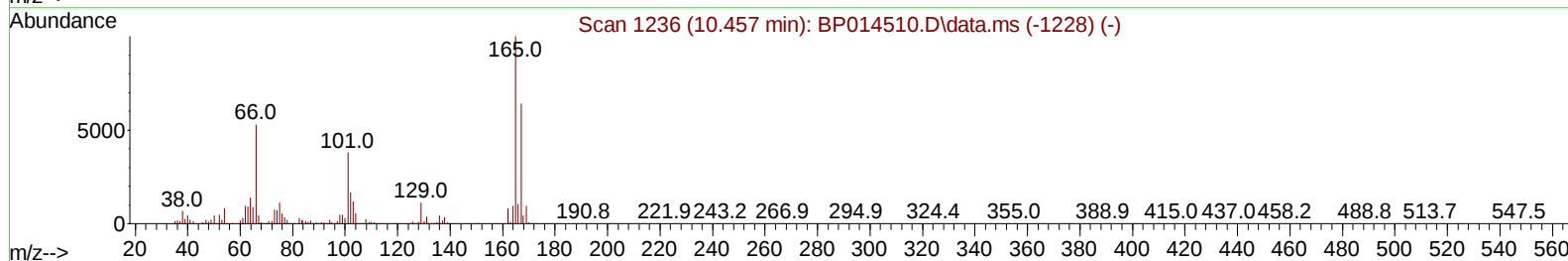
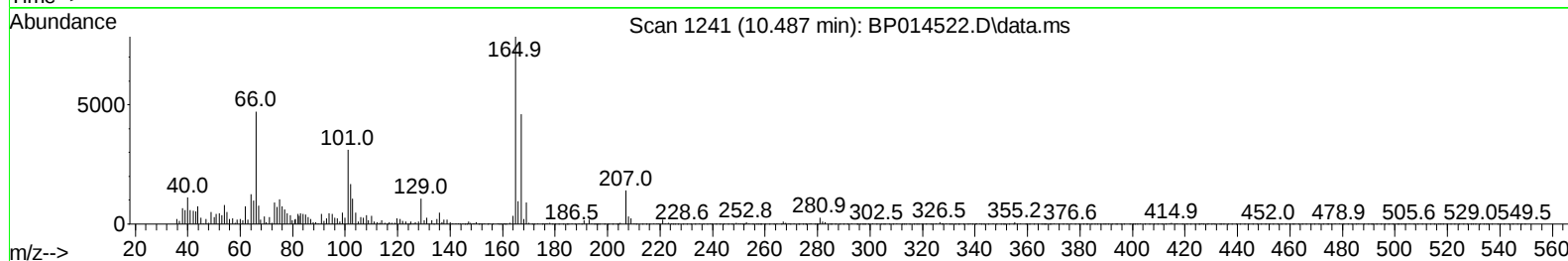
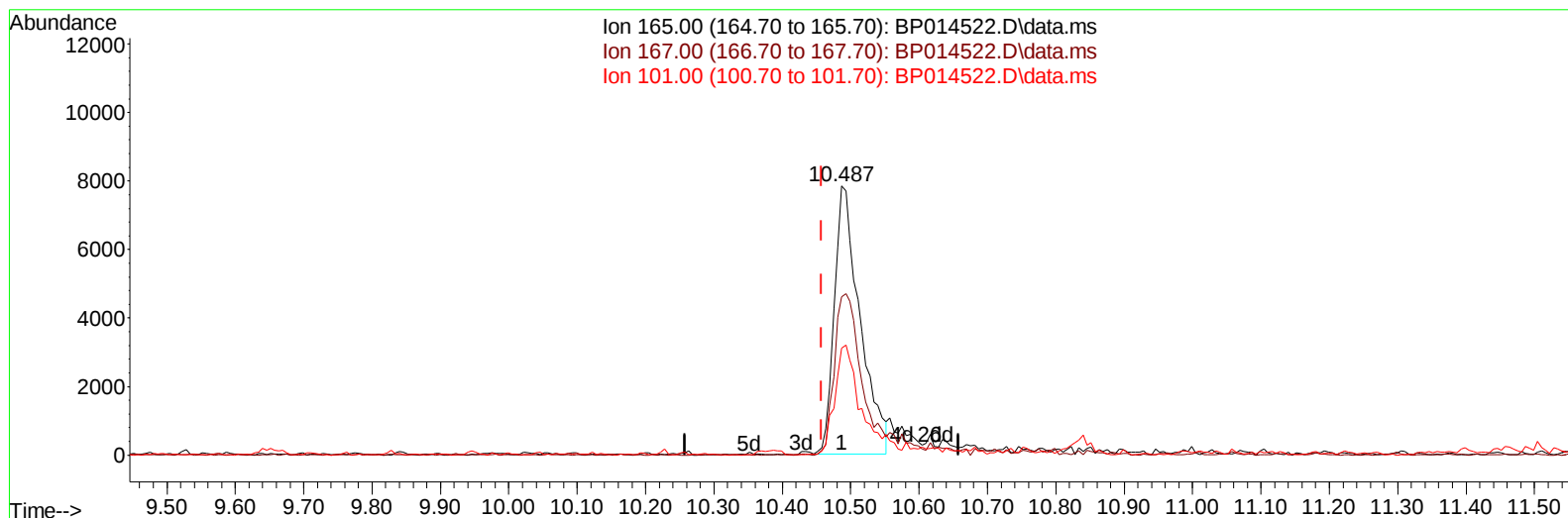
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TIC: BP014522.D\data.ms

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

10.487min (+ 0.029) 2.01 ng/ul

response 20481

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.80	58.80
101.00	39.20	39.82
0.00	0.00	0.00

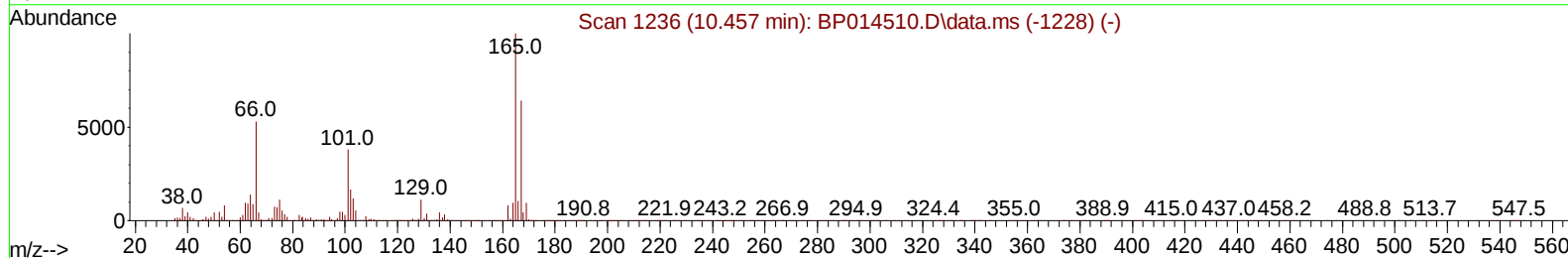
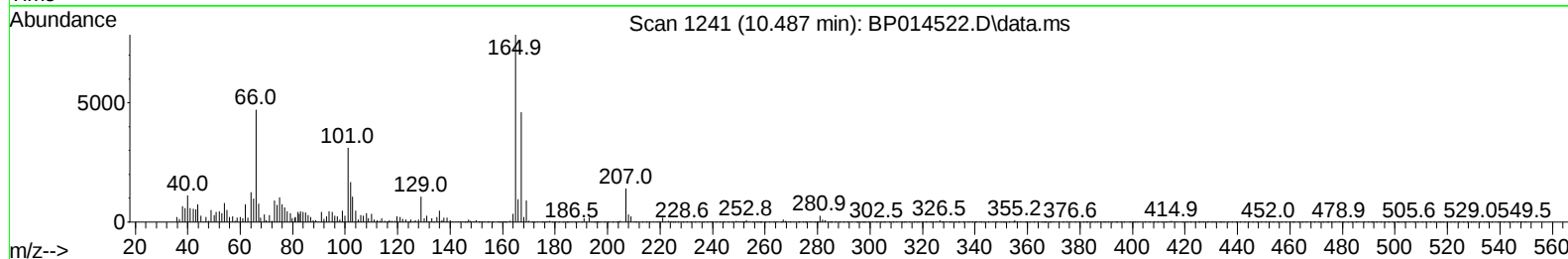
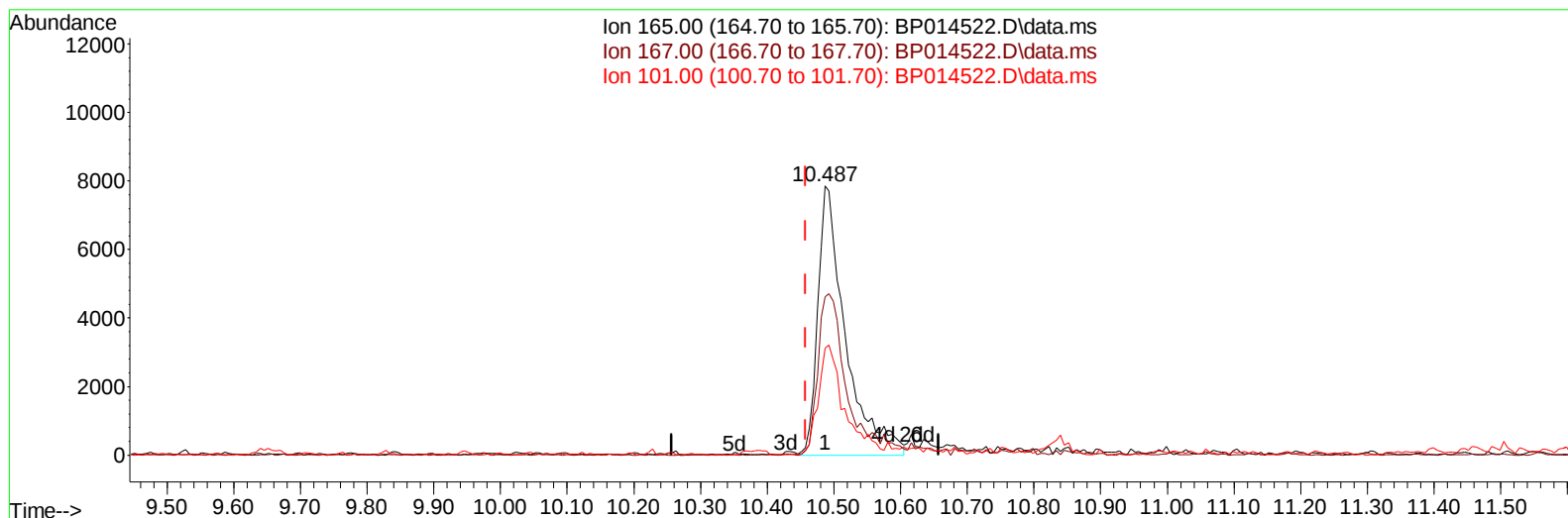
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Instrument :
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TIC: BP014522.D\data.ms

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

10.487min (+ 0.029) 2.22 ng/ul m

response 22606

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.80	58.80
101.00	39.20	39.82
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	8.016	152	143997	20.000	ng/ul	0.01
20) Naphthalene-d8	10.840	136	605515	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.663	164	358383	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	650332	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	573640	20.000	ng/ul	0.00
88) Perylene-d12	24.027	264	719787	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	2266	0.610	ng/uL	0.01
4) Pyridine-d5	3.864	84	15879m	1.669	ng/ul	0.05
7) Phenol-d5	7.211	99	10264	0.773	ng/ul	0.04
9) Bis-(2-Chloroethyl)eth...	7.346	67	24760	2.906	ng/ul	0.01
11) 2-Chlorophenol-d4	7.558	132	20717	2.158	ng/ul	0.02
15) 4-Methylphenol-d8	8.752	113	19353	1.807	ng/ul	0.02
21) Nitrobenzene-d5	9.199	128	13551	2.772	ng/ul	0.01
24) 2-Nitrophenol-d4	9.928	143	12388	2.213	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.487	165	22606m	2.218	ng/ul	0.03
31) 4-Chloroaniline-d4	11.005	131	29934	2.229	ng/ul	0.02
46) Dimethylphthalate-d6	14.075	166	85812	2.953	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	93519	2.900	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.663	176	71601	2.955	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.528	188	104570	3.323	ng/ul	0.00
81) Pyrene-d10	19.757	212	111549	2.895	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.869	264	118709	3.104	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.423	88	105344	25.816	ng/uL	94

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

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