

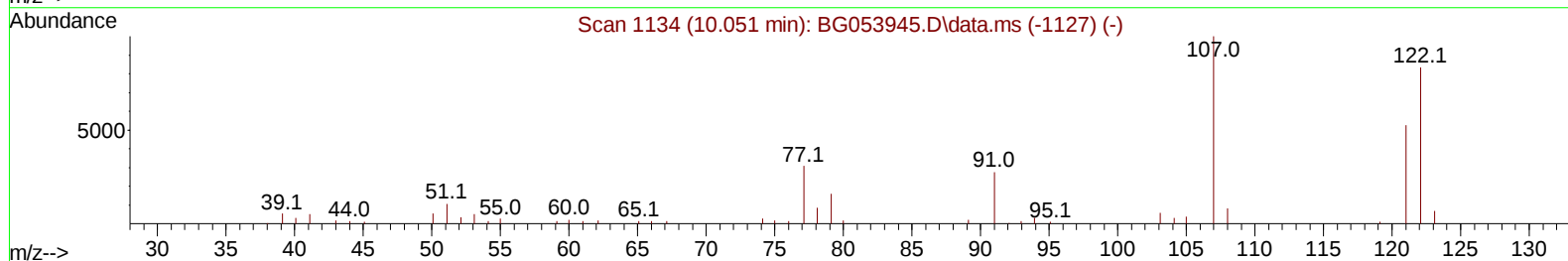
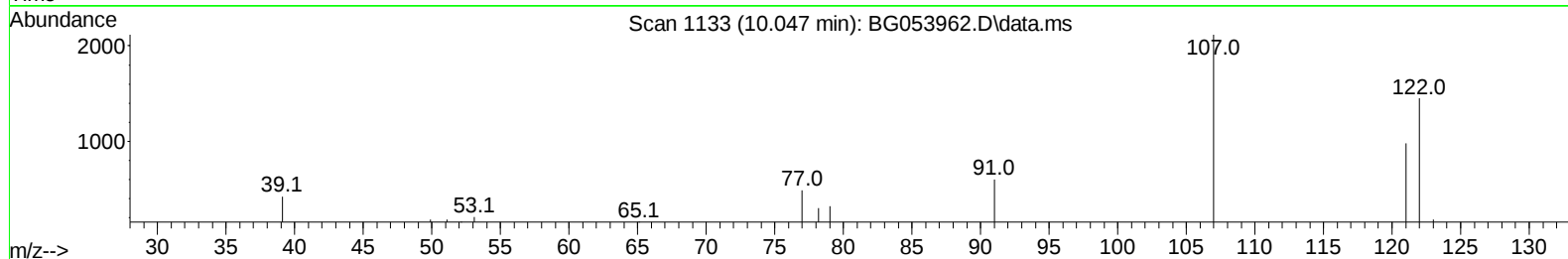
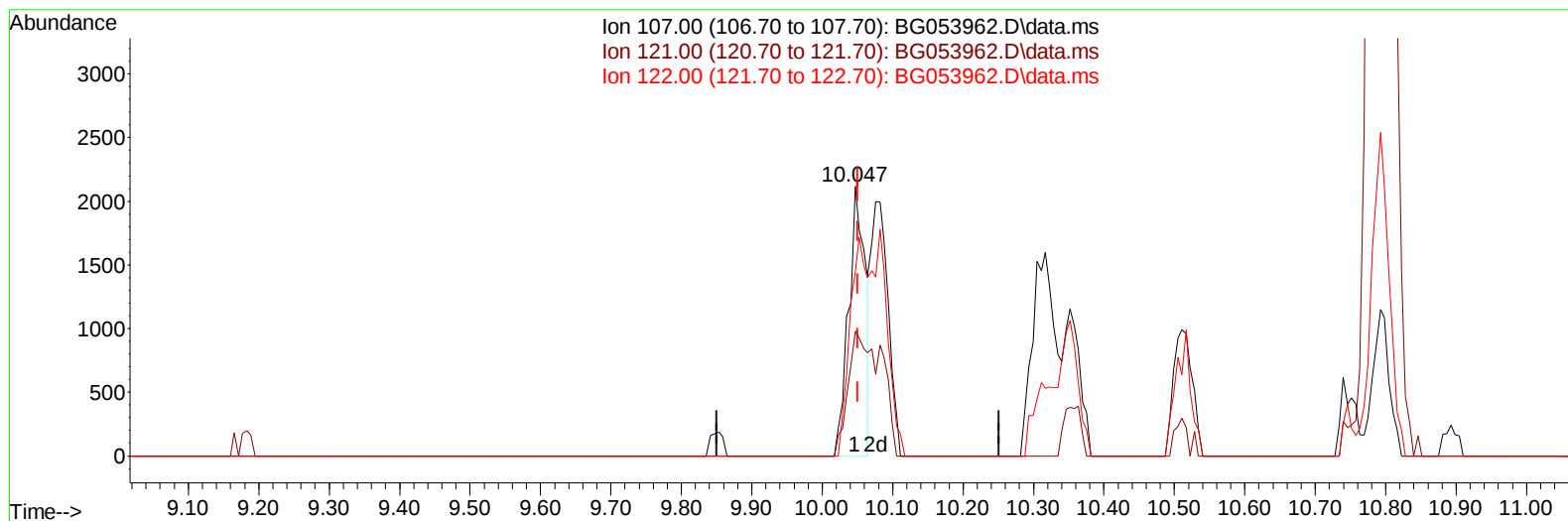
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062022\
 Data File : BG053962.D
 Acq On : 20 Jun 2022 23:50
 Operator : CG/JU
 Sample : N3358-15
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW4S8

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 06/21/2022
 Supervised By :mohammad ahmed 06/27/2022

Quant Time: Jun 21 00:32:46 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 20 15:29:13 2022
 Response via : Initial Calibration



TIC: BG053962.D\data.ms

(26) 2,4-Dimethylphenol

10.047min (-0.005) 2.64 ng/u1

response 3479

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	49.90	46.36
122.00	83.40	68.53
0.00	0.00	0.00

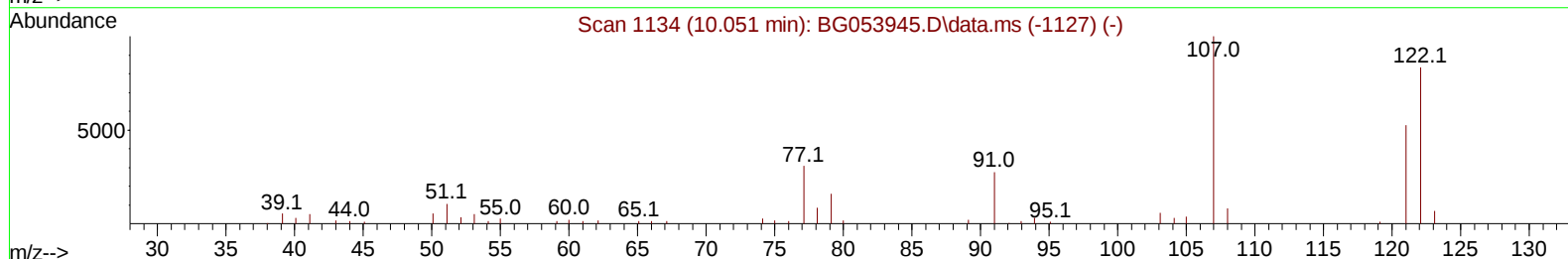
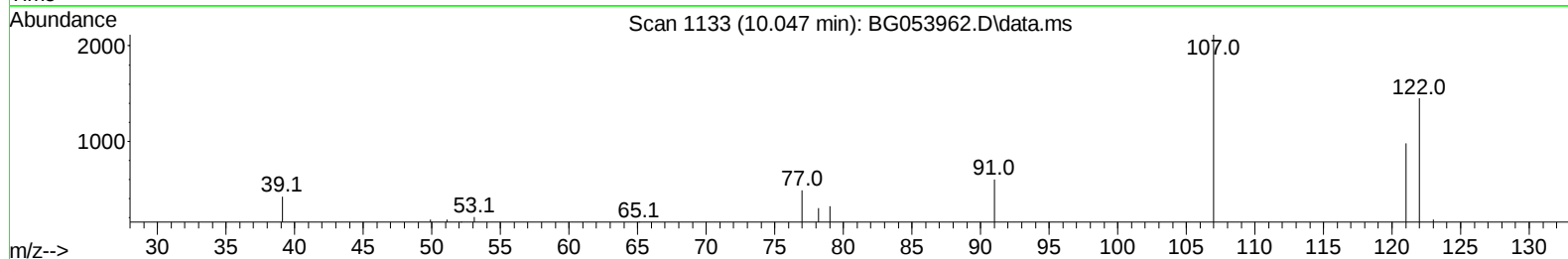
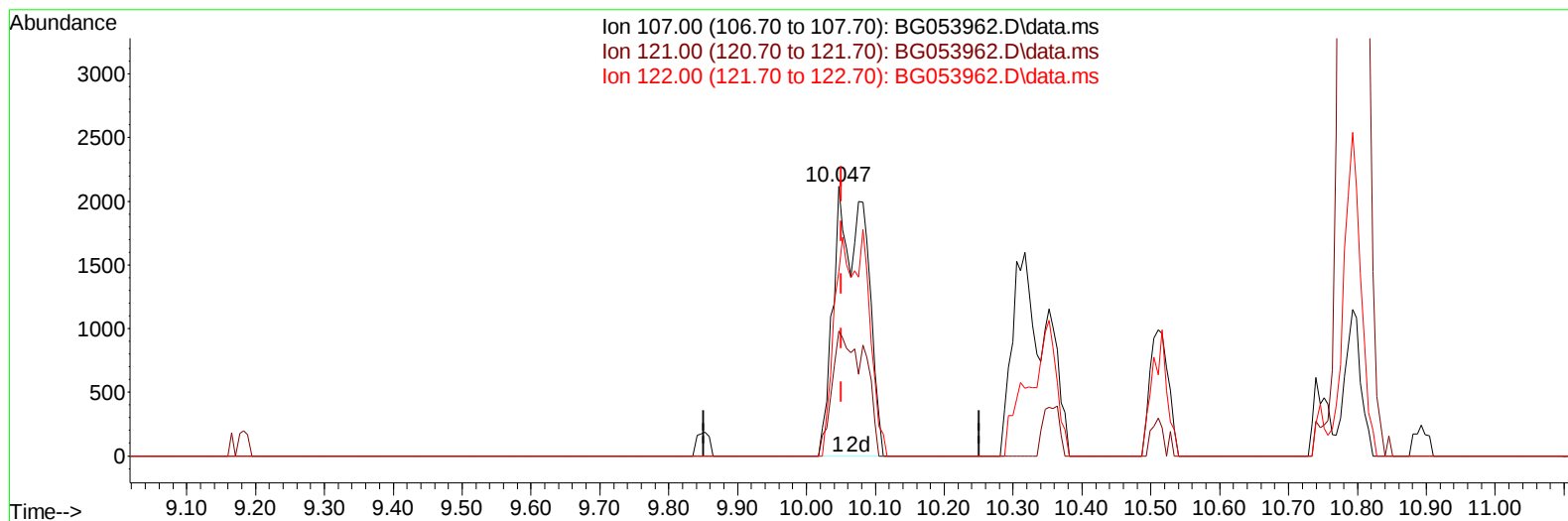
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TIC: BG053962.D\data.ms

(26) 2,4-Dimethylphenol

10.047min (-0.005) 5.18 ng/u1 m

response 6842

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	49.90	46.36
122.00	83.40	68.53
0.00	0.00	0.00

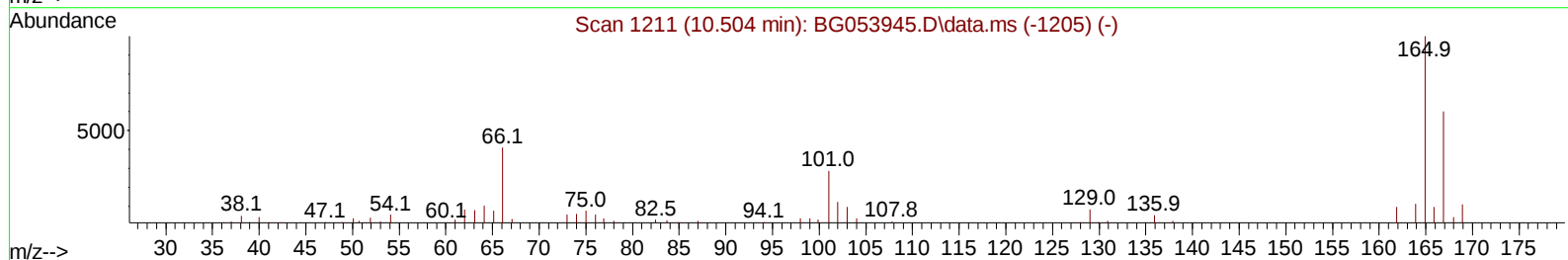
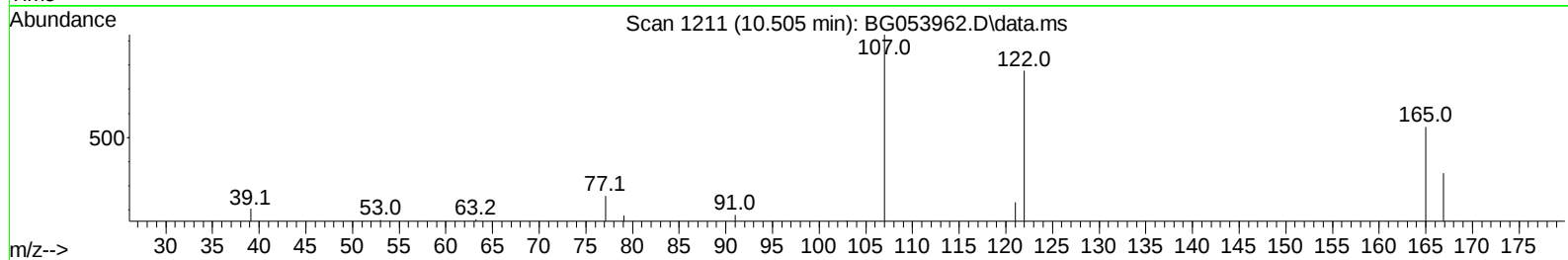
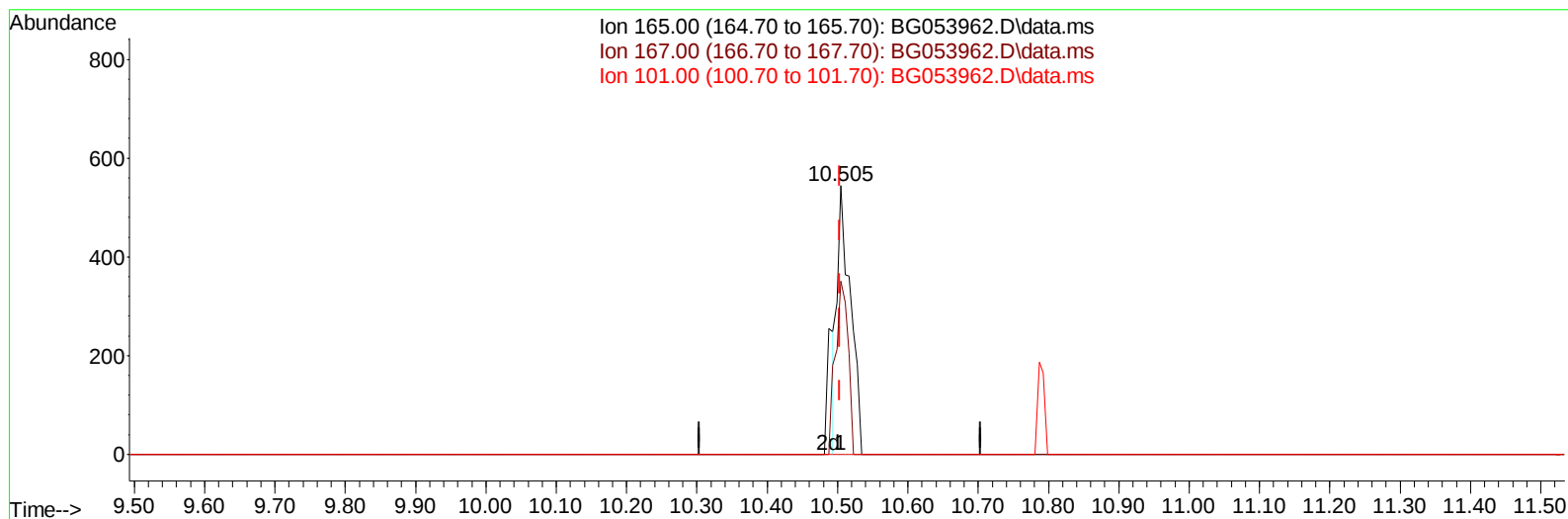
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TIC: BG053962.D\data.ms

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

10.505min (+ 0.001) 0.56 ng/ul

response 708

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	61.80	64.52
101.00	32.90	0.00#
0.00	0.00	0.00

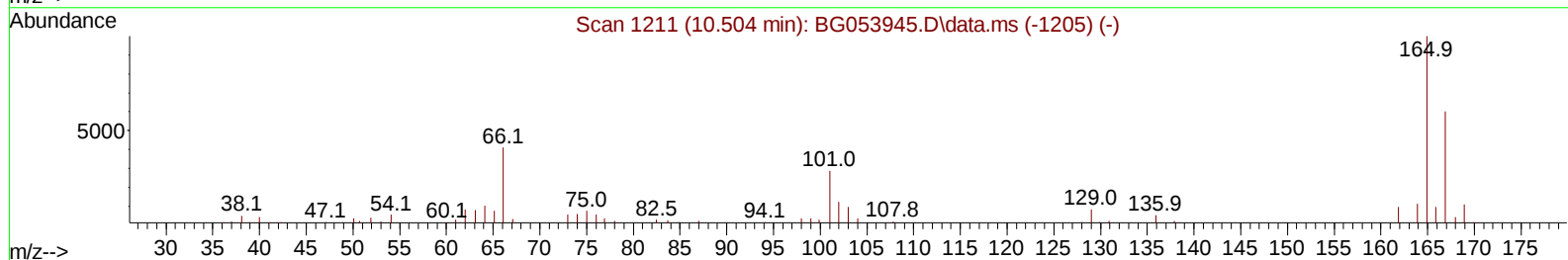
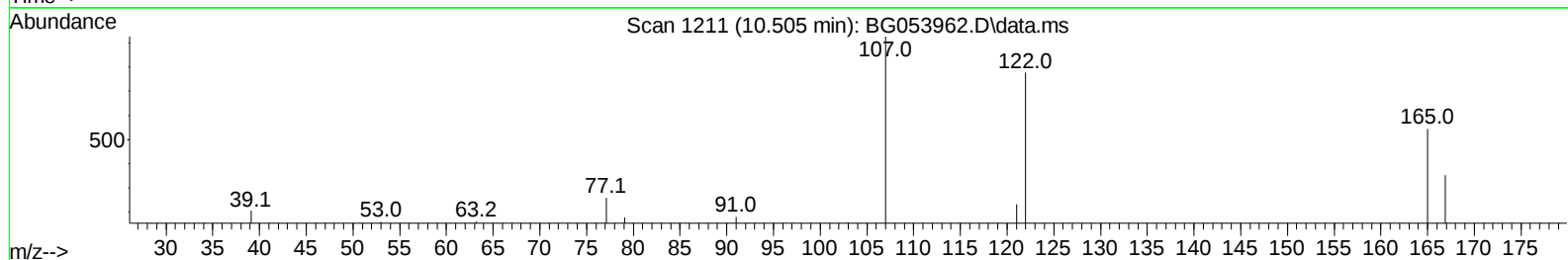
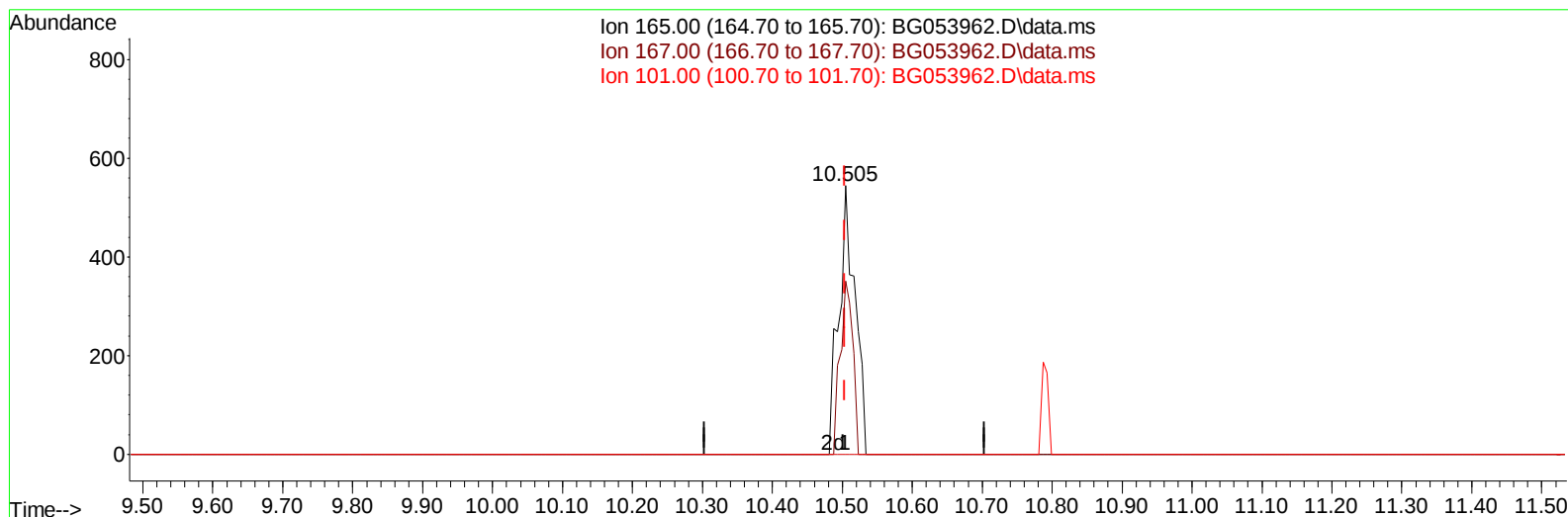
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TIC: BG053962.D\data.ms

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

10.505min (+ 0.001) 0.69 ng/ul m

response 886

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	61.80	64.52
101.00	32.90	0.00#
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	8.078	152	18685	20.000	ng/ul	0.00
20) Naphthalene-d8	10.887	136	73094	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.700	164	56934	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.444	188	155786	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.721	240	178582	20.000	ng/ul	# 0.00
88) Perylene-d12	24.976	264	186036	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.507	96	1596	4.064	ng/uL	0.00
4) Pyridine-d5	3.918	84	13555	12.690	ng/ul	0.00
7) Phenol-d5	7.232	99	2533	1.892	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.397	67	18421	22.813	ng/ul	0.00
11) 2-Chlorophenol-d4	7.608	132	908	0.851	ng/ul	0.00
15) 4-Methylphenol-d8	8.777	113	9188	8.050	ng/ul	0.00
21) Nitrobenzene-d5	9.236	128	13090	24.127	ng/ul	0.00
24) 2-Nitrophenol-d4	9.958	143	348	0.604	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.505	165	886m	0.695	ng/ul	0.00
31) 4-Chloroaniline-d4	11.016	131	36781	22.828	ng/ul	0.00
46) Dimethylphthalate-d6	14.101	166	120678	26.943	ng/ul	0.00
49) Acenaphthylene-d8	14.394	160	121976	25.045	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0	0.000	ng/ul	
60) Fluorene-d10	15.693	176	104190	25.897	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0	0.000	ng/ul	
73) Anthracene-d10	17.544	188	194989	27.715	ng/ul	0.00
81) Pyrene-d10	19.829	212	265143	28.476	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.753	264	262121	27.913	ng/ul	0.00
Target Compounds						
8) Phenol	7.262	94	16055	11.495	ng/ul	Qvalue 97
26) 2,4-Dimethylphenol	10.047	107	6842m	5.184	ng/ul	

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

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