

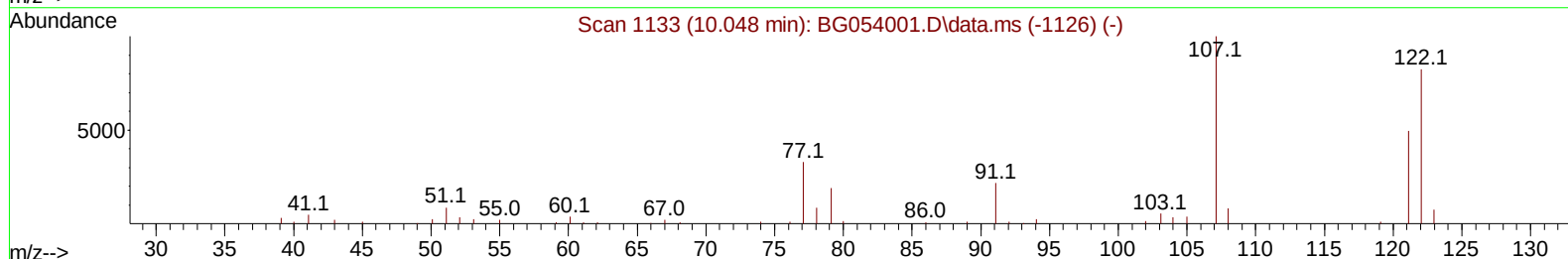
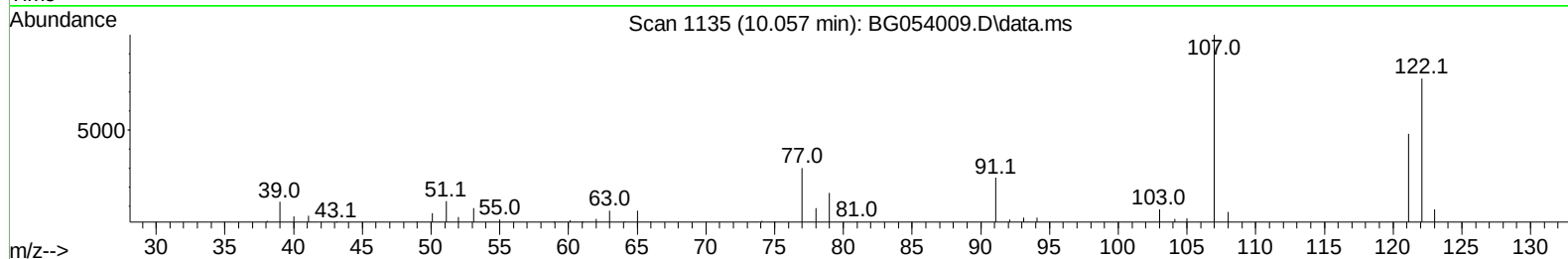
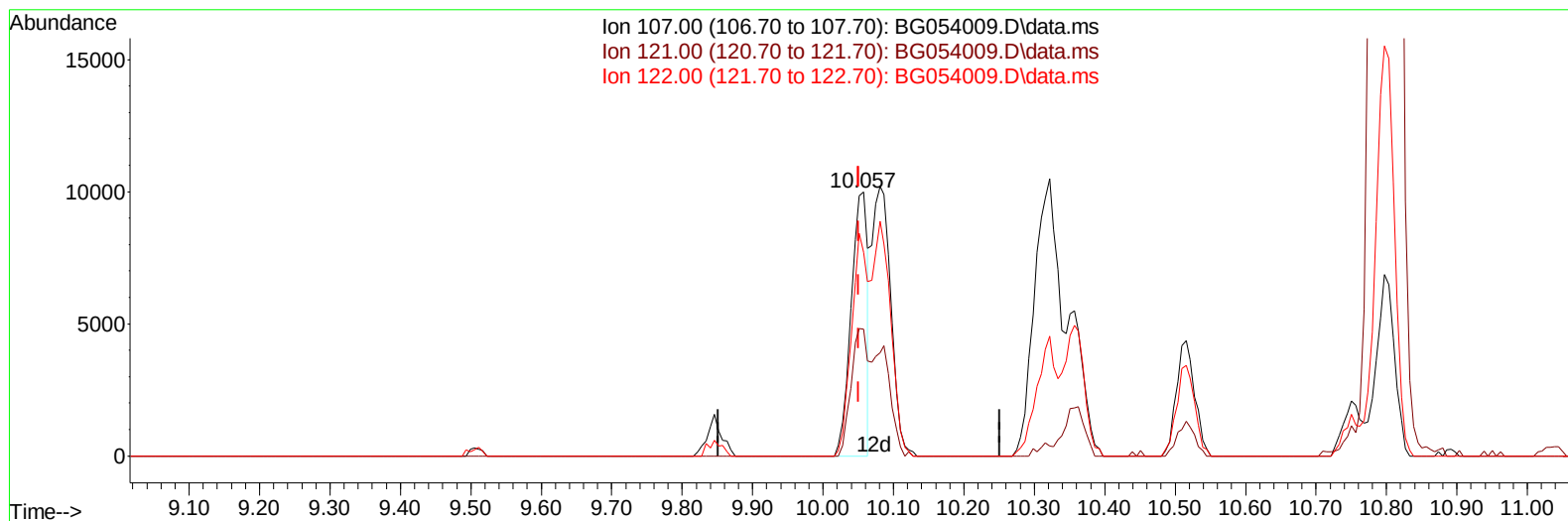
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062122\
 Data File : BG054009.D
 Acq On : 22 Jun 2022 13:34
 Operator : CG/JU
 Sample : N3358-18DL2 50X
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW4T1DL2

Manual IntegrationsAPPROVED

Quant Time: Jun 23 01:13:50 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

Reviewed By :Jagrut Upadhyay 06/23/2022
 Supervised By :mohammad ahmed 06/28/2022



TIC: BG054009.D\data.ms

(26) 2,4-Dimethylphenol

10.057min (+ 0.006) 9.23 ng/ul

response 16179

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	49.90	48.08
122.00	83.40	77.08
0.00	0.00	0.00

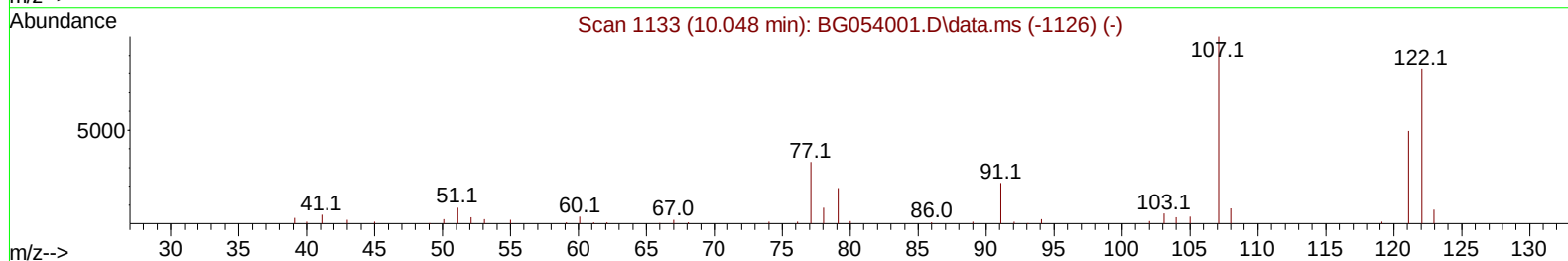
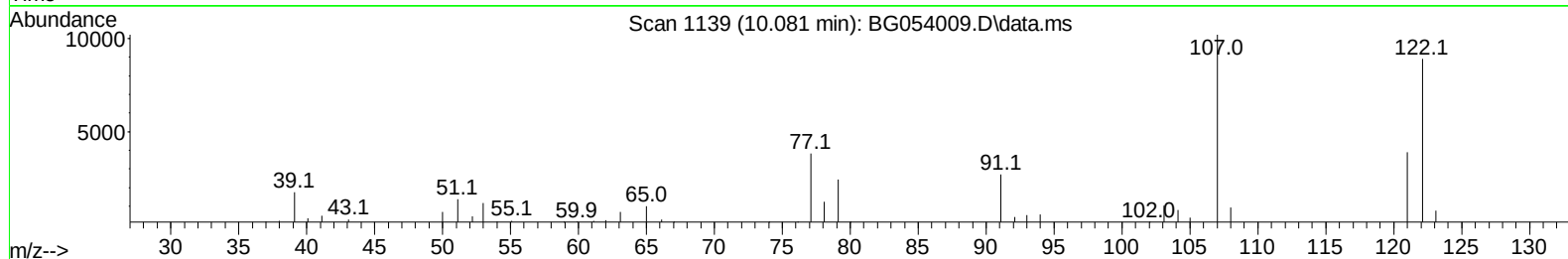
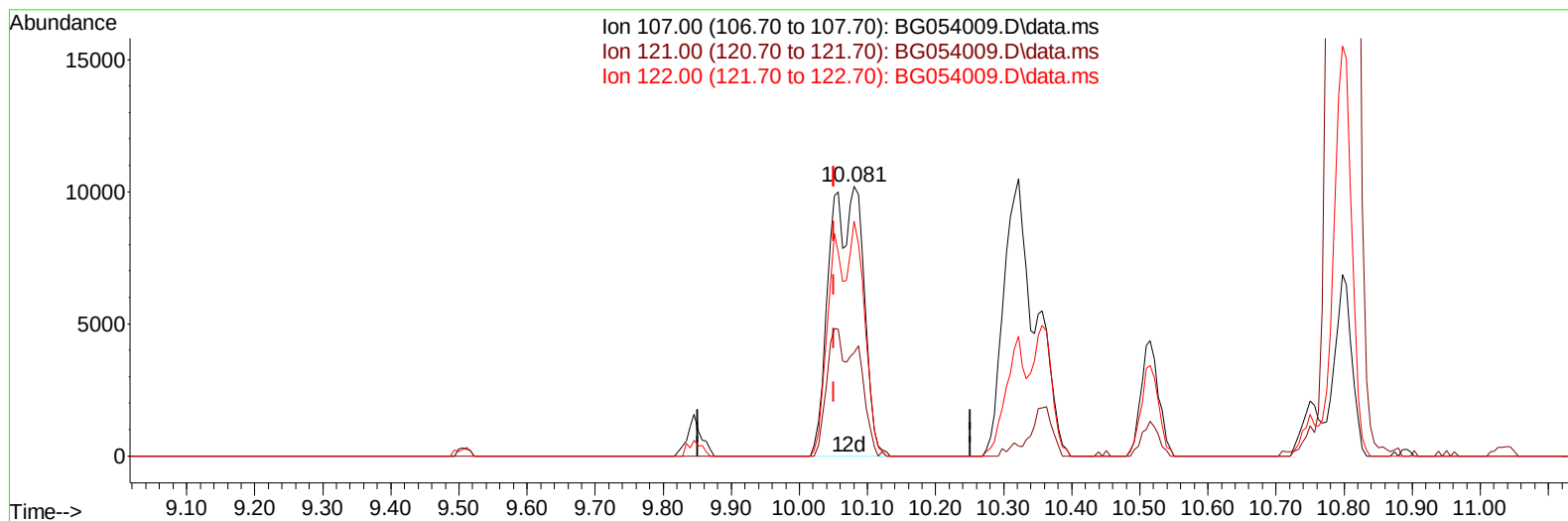
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(26) 2,4-Dimethylphenol

10.081min (+ 0.030) 20.16 ng/ul m

response 35357

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	49.90	38.40#
122.00	83.40	87.17
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	8.077	152	25137	20.000	ng/ul	0.00
20) Naphthalene-d8	10.891	136	97121	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.705	164	75504	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.448	188	196581	20.000	ng/ul	0.00
79) Chrysene-d12	21.720	240	219467	20.000	ng/ul	# 0.00
88) Perylene-d12	24.981	264	221957	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0	0.000	ng/uL	
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	0.000	99	0d	0.000	ng/ul	
9) Bis-(2-Chloroethyl)eth...	7.396	67	552	0.508	ng/ul	0.00
11) 2-Chlorophenol-d4	7.607	132	416	0.290	ng/ul	0.00
15) 4-Methylphenol-d8	8.776	113	354	0.231	ng/ul	0.00
21) Nitrobenzene-d5	9.240	128	266	0.369	ng/ul	0.00
24) 2-Nitrophenol-d4	9.969	143	229	0.299	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.504	165	734	0.433	ng/ul	0.00
31) 4-Chloroaniline-d4	11.021	131	709	0.331	ng/ul	0.00
46) Dimethylphthalate-d6	14.105	166	2800	0.471	ng/ul	0.00
49) Acenaphthylene-d8	14.399	160	3501	0.542	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0	0.000	ng/ul	
60) Fluorene-d10	15.692	176	3045	0.571	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.548	188	5589	0.630	ng/ul	0.00
81) Pyrene-d10	19.828	212	6624	0.579	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.752	264	6195	0.553	ng/ul	0.00
Target Compounds						
						Qvalue
8) Phenol	7.260	94	172808	91.966	ng/ul	94
18) 4-Methylphenol	8.847	108	8165	5.233	ng/ul	83
26) 2,4-Dimethylphenol	10.081	107	35357m	20.160	ng/ul	

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

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