

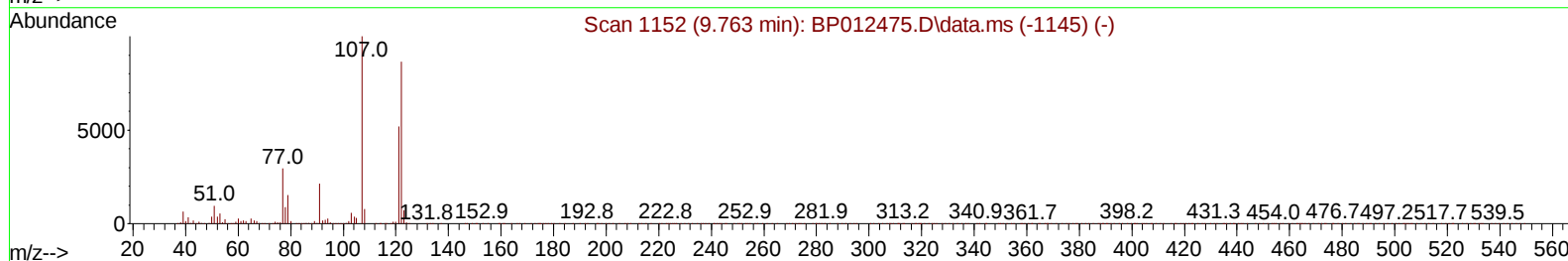
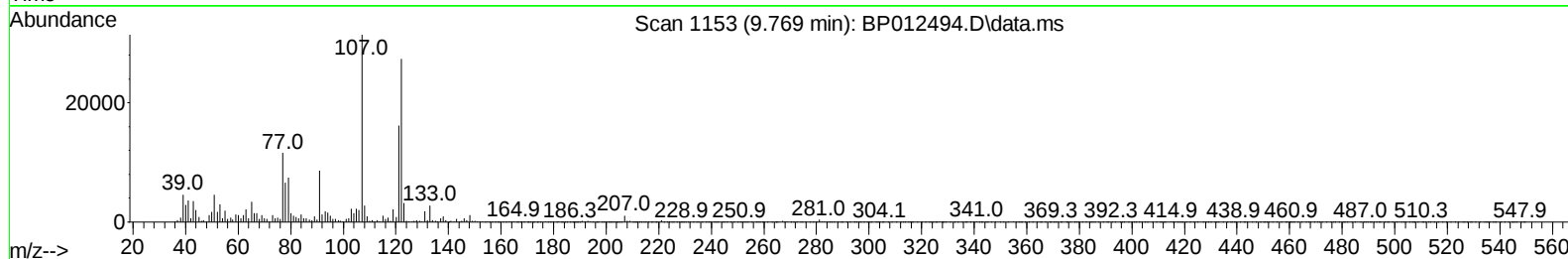
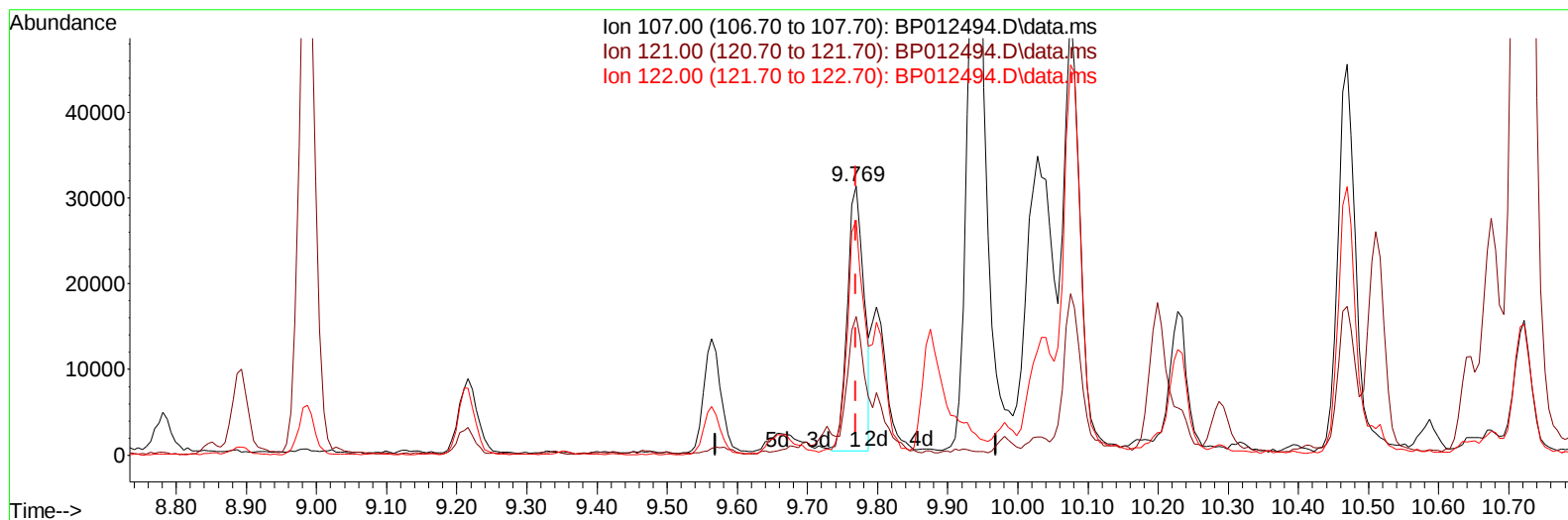
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012494.D
 Acq On : 03 Nov 2022 17:48
 Operator : CG/JU
 Sample : N5319-09
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHAA2

Manual IntegrationsAPPROVED

Quant Time: Nov 03 23:05:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 03 00:09:04 2022
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/05/2022
 Supervised By :mohammad ahmed 11/06/2022



TIC: BP012494.D\data.ms

(26) 2,4-Dimethylphenol

9.769min (+ 0.000) 2.44 ng/u1

response 51346

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	51.10	51.64
122.00	88.20	87.29
0.00	0.00	0.00

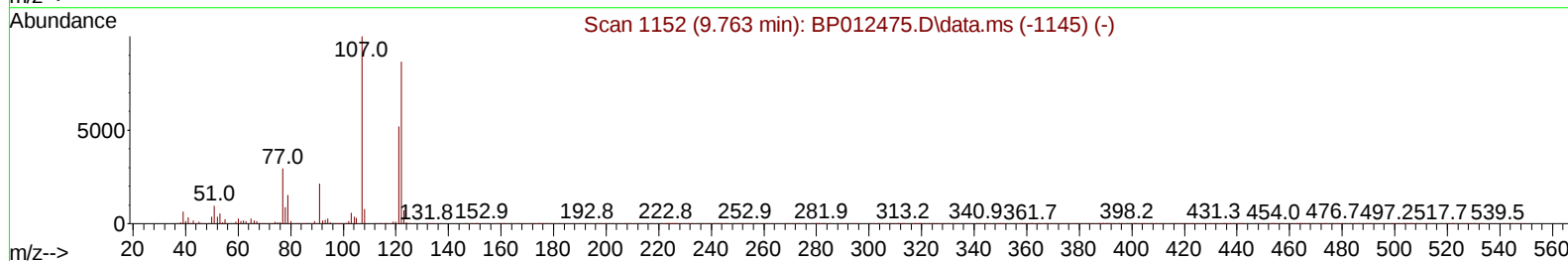
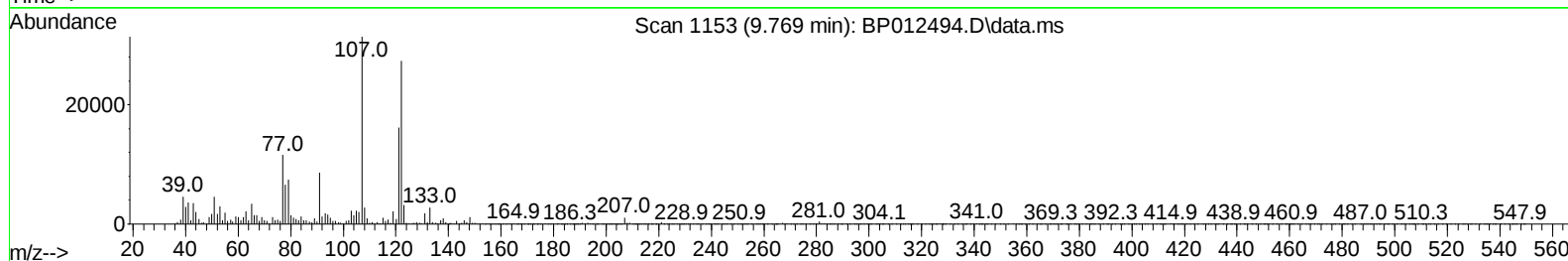
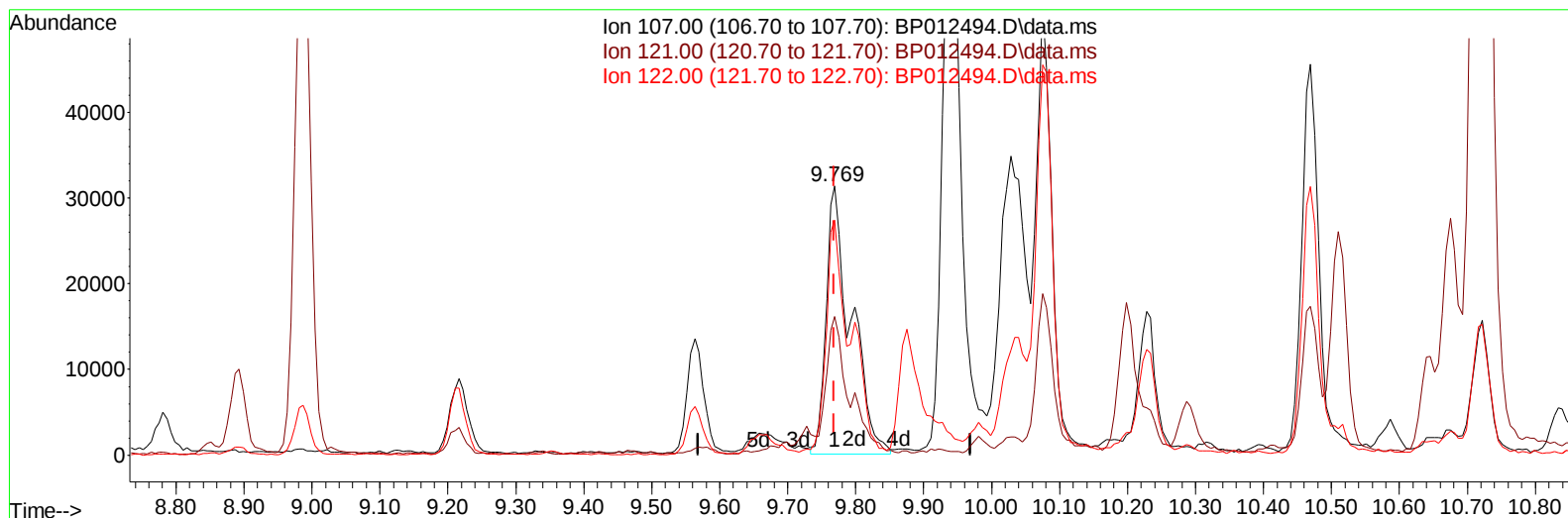
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TIC: BP012494.D\data.ms

(26) 2,4-Dimethylphenol

9.769min (+ 0.000) 3.75 ng/u1 m

response 79024

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	51.10	51.64
122.00	88.20	87.29
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	7.781	152	247213	20.000	ng/ul	0.00
20) Naphthalene-d8	10.581	136	1189625	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.434	164	816063	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	2166754	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.292	240	2274265	20.000	ng/ul	0.00
88) Perylene-d12	23.674	264	2089885	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	25148	4.096	ng/uL	0.00
4) Pyridine-d5	3.658	84	178589	10.126	ng/ul	0.00
7) Phenol-d5	6.963	99	166280	7.499	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	422769	31.939	ng/ul	0.00
11) 2-Chlorophenol-d4	7.316	132	436630	26.965	ng/ul	0.00
15) 4-Methylphenol-d8	8.505	113	314900	17.956	ng/ul	0.00
21) Nitrobenzene-d5	8.952	128	297858	38.610	ng/ul	0.00
24) 2-Nitrophenol-d4	9.669	143	319530	39.906	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	574880	33.129	ng/ul	0.00
31) 4-Chloroaniline-d4	10.734	131	712995	27.595	ng/ul	0.00
46) Dimethylphthalate-d6	13.851	166	2000521	35.811	ng/ul	0.00
49) Acenaphthylene-d8	14.128	160	2170920	34.005	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	66988	6.336	ng/ul	0.02
60) Fluorene-d10	15.434	176	1668566	34.886	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	360267	31.936	ng/ul	0.00
73) Anthracene-d10	17.298	188	3337678	35.750	ng/ul	0.00
81) Pyrene-d10	19.539	212	3997293	35.008	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.521	264	3703258	36.266	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.269	88	7161	1.097	ng/ul#	92
8) Phenol	6.993	94	197696	8.575	ng/ul	95
10) Bis(2-Chloroethyl)ether	7.216	93	54664	2.954	ng/ul	95
12) 2-Chlorophenol	7.352	128	23698	1.344	ng/ul	98
13) 2-Methylphenol	8.240	108	81410	4.592	ng/ul	99
16) Acetophenone	8.616	105	36725	1.348	ng/ul	90
18) 4-Methylphenol	8.563	108	283316	14.620	ng/ul	95
26) 2,4-Dimethylphenol	9.769	107	79024m	3.748	ng/ul	
27) Bis(2-Chloroethoxy)met...	10.004	93	103373	3.835	ng/ul#	34
30) Naphthalene	10.634	128	281466	4.464	ng/ul	100
59) Diethylphthalate	15.263	149	168169	2.936	ng/ul	99
67) N-Nitrosodiphenylamine	15.704	169	121934	1.998	ng/ul	99

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

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