

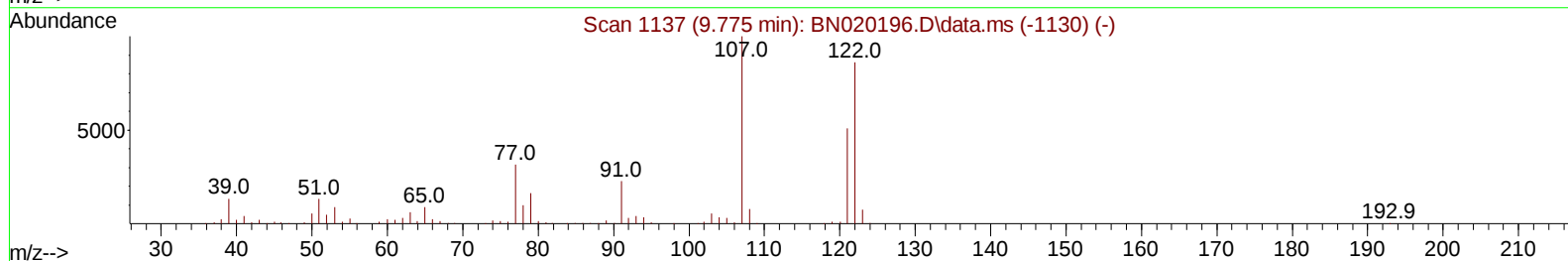
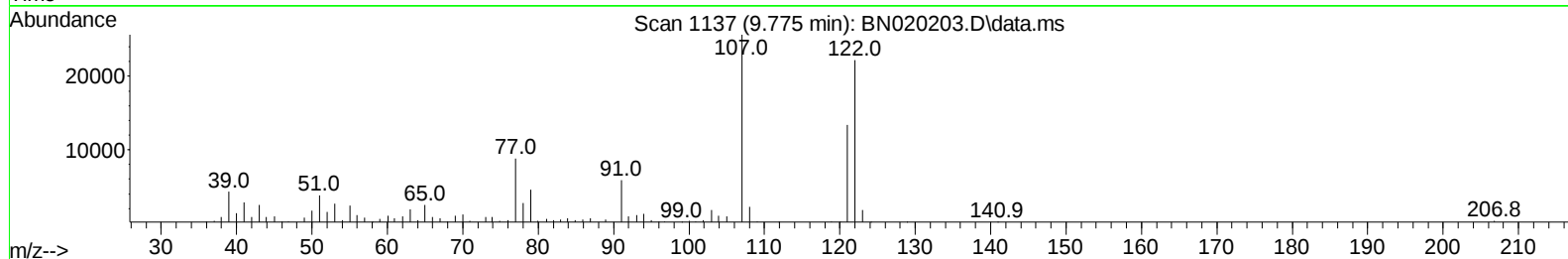
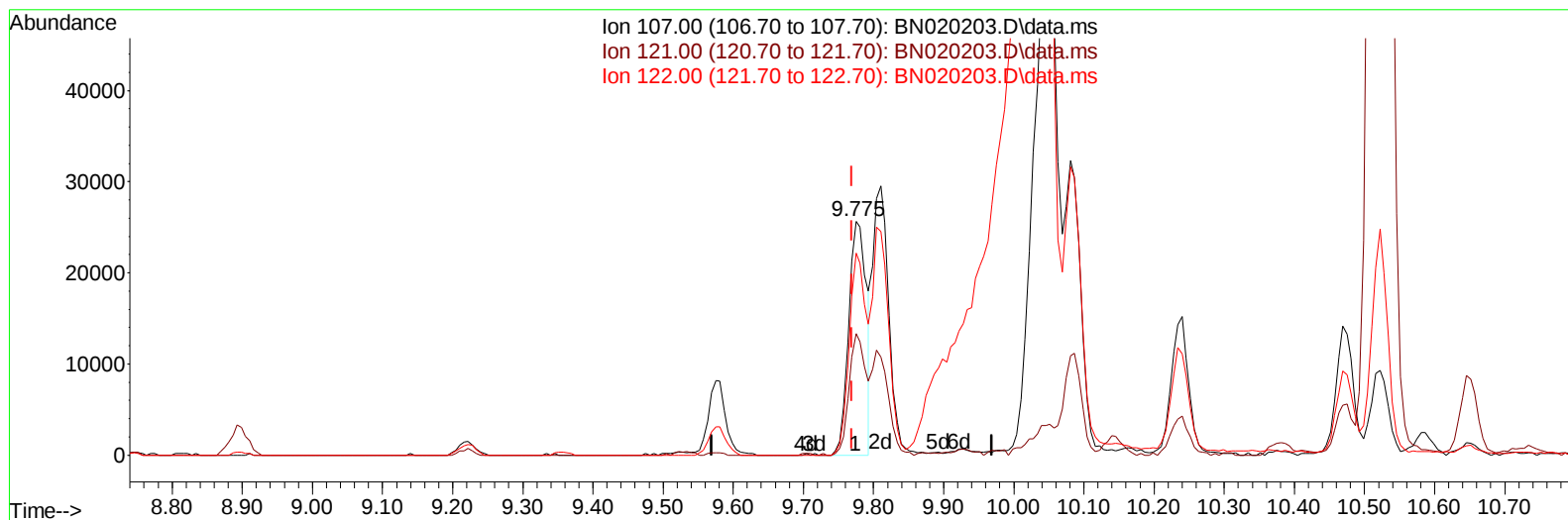
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061522\
 Data File : BN020203.D
 Acq On : 16 Jun 2022 14:18
 Operator : CG/JU
 Sample : N3285-01DL 5X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EW4P5DL

Manual IntegrationsAPPROVED

Quant Time: Jun 17 02:27:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 16 00:50:14 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/17/2022
 Supervised By :mohammad ahmed 06/22/2022



TIC: BN020203.D\data.ms

(26) 2,4-Dimethylphenol

9.775min (+ 0.006) 3.07 ng/u1

response 45637

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	47.90	52.03
122.00	82.90	86.37
0.00	0.00	0.00

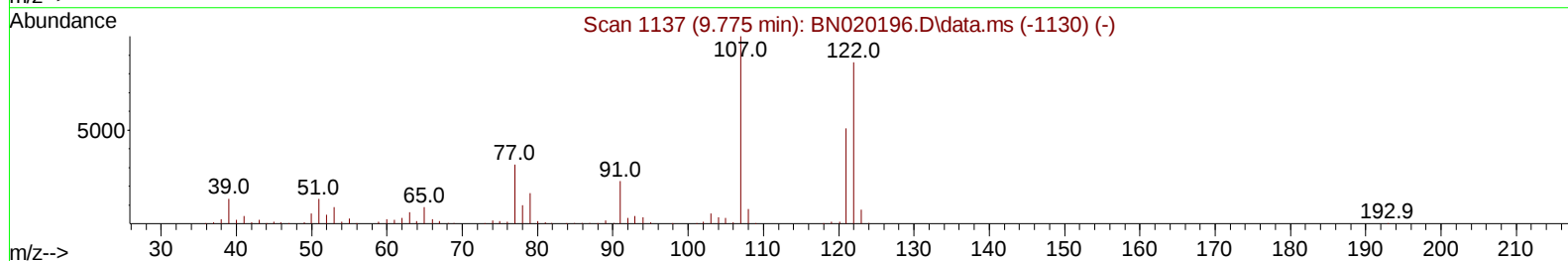
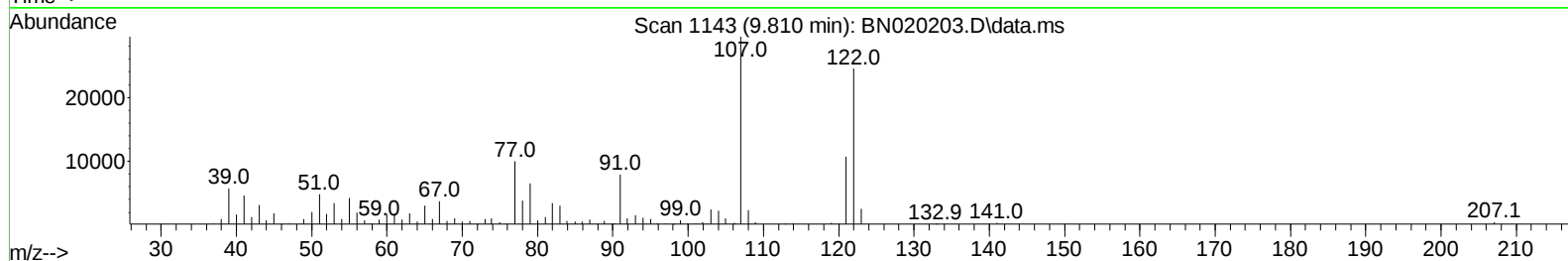
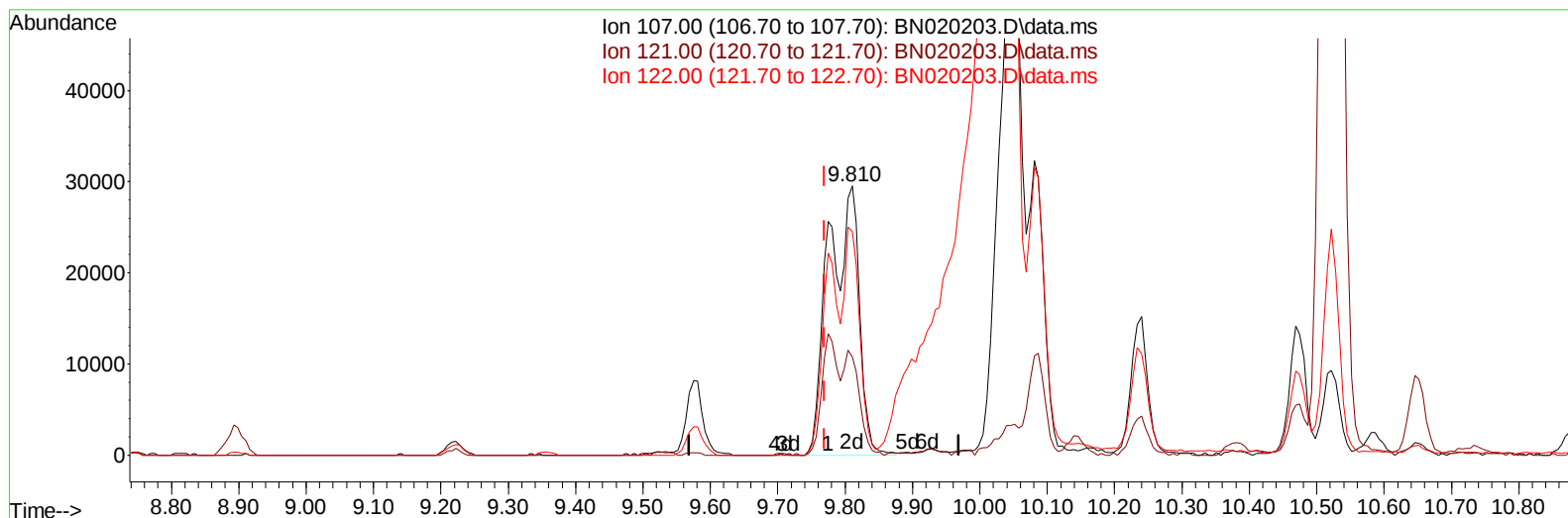
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(26) 2,4-Dimethylphenol

9.810min (+ 0.041) 6.26 ng/ul m

response 92861

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	47.90	36.34#
122.00	82.90	83.04
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	7.793	152	172273	20.000	ng/ul	0.00
20) Naphthalene-d8	10.581	136	803061	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.422	164	519385	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.163	188	1061993	20.000	ng/ul	0.00
79) Chrysene-d12	21.357	240	869886	20.000	ng/ul	0.00
88) Perylene-d12	23.674	264	703051	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	3962	0.996	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.987	99	22599	1.585	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.128	67	44055	4.946	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	49385	4.264	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	42003	3.406	ng/ul	0.01
21) Nitrobenzene-d5	8.946	128	29917	4.897	ng/ul	0.00
24) 2-Nitrophenol-d4	9.669	143	31539	4.864	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	57520	5.082	ng/ul	0.01
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	13.839	166	206257	5.458	ng/ul	0.00
49) Acenaphthylene-d8	14.116	160	206426	4.710	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.416	176	172269	5.509	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.534	200	28419	4.620	ng/ul	0.00
73) Anthracene-d10	17.263	188	262362	5.661	ng/ul	0.00
81) Pyrene-d10	19.557	212	282920	5.999	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.521	264	185829	5.402	ng/ul	-0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.293	88	7203	1.715	ng/uL#	88
8) Phenol	7.016	94	1359966	89.593	ng/ul	99
18) 4-Methylphenol	8.587	108	138434	10.748	ng/ul	94
26) 2,4-Dimethylphenol	9.810	107	92861m	6.256	ng/ul	

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

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