

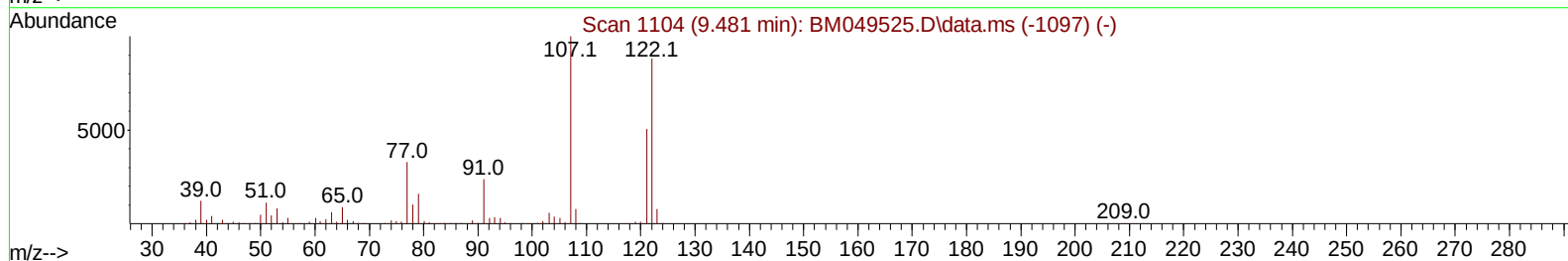
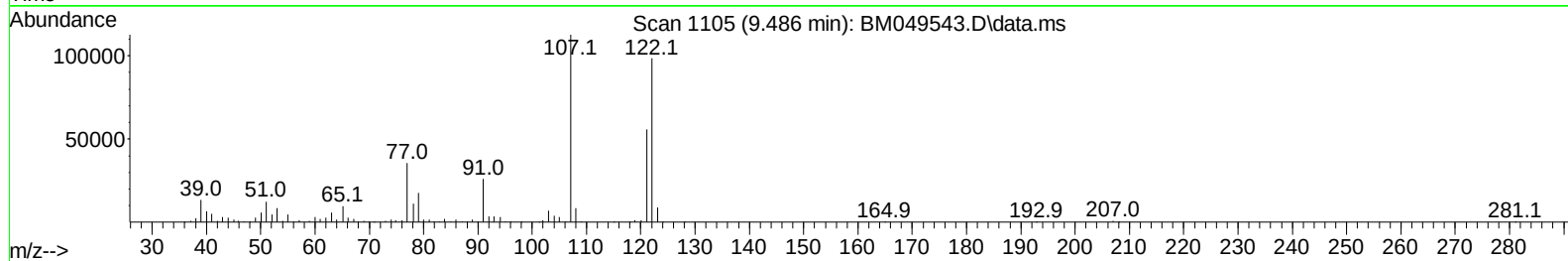
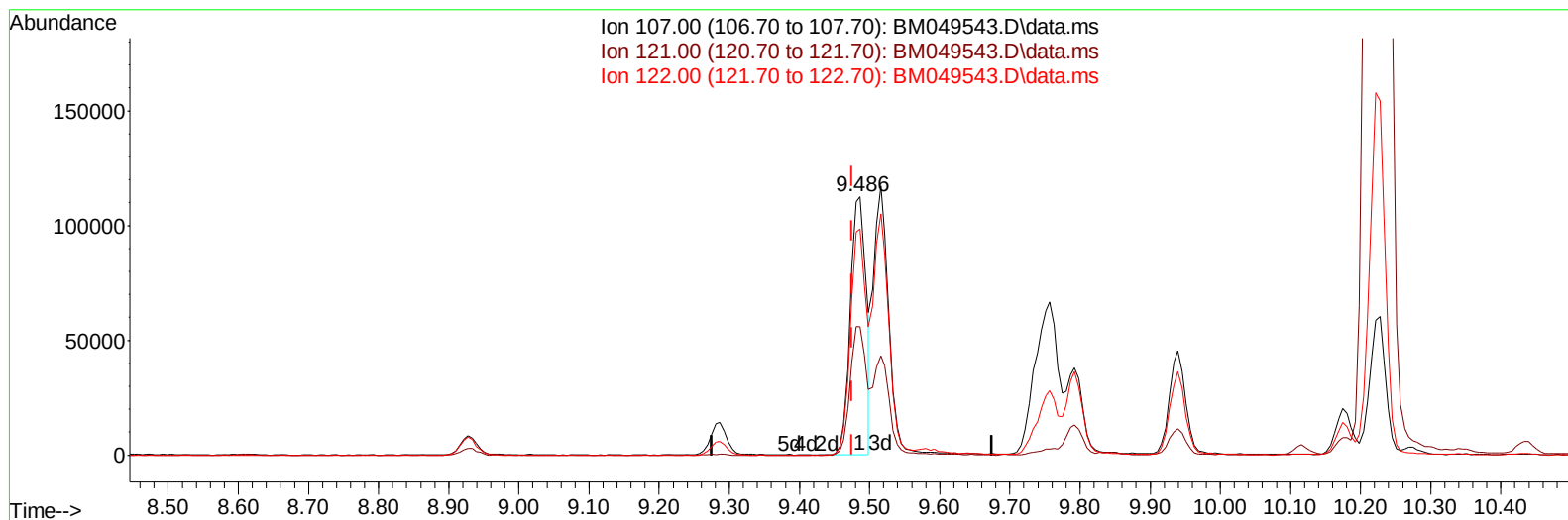
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013025\
Data File : BM049543.D
Acq On : 31 Jan 2025 01:42
Operator : RC/JU
Sample : Q1184-15DL 200X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E2966DL

Manual IntegrationsAPPROVED

Quant Time: Jan 31 08:52:56 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM012825.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jan 28 14:45:18 2025
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/31/2025
Supervised By :mohammad ahmed 02/04/2025



TIC: BM049543.D\data.ms

(26) 2,4-Dimethylphenol

9.486min (+ 0.012) 9.37 ng/u1

response 178279

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	49.50	49.59
122.00	85.90	87.58
0.00	0.00	0.00

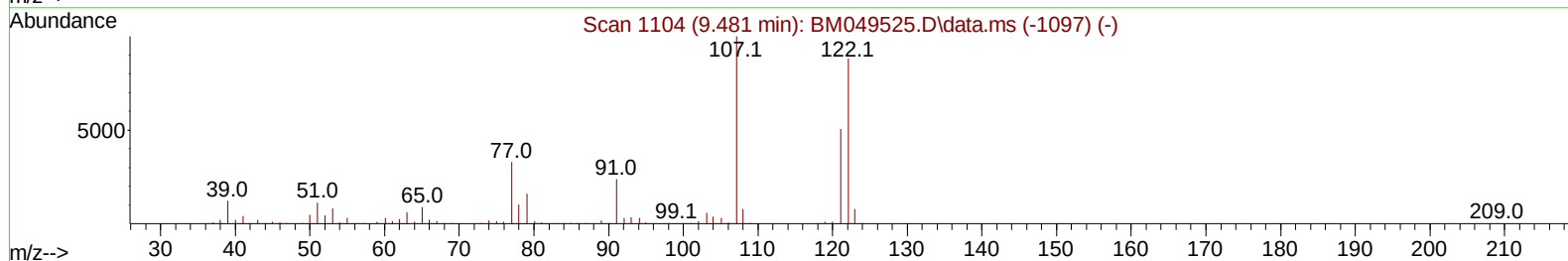
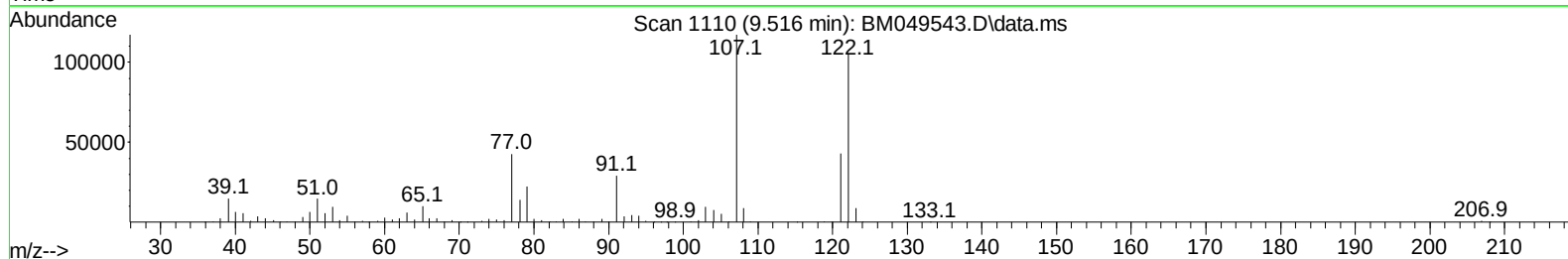
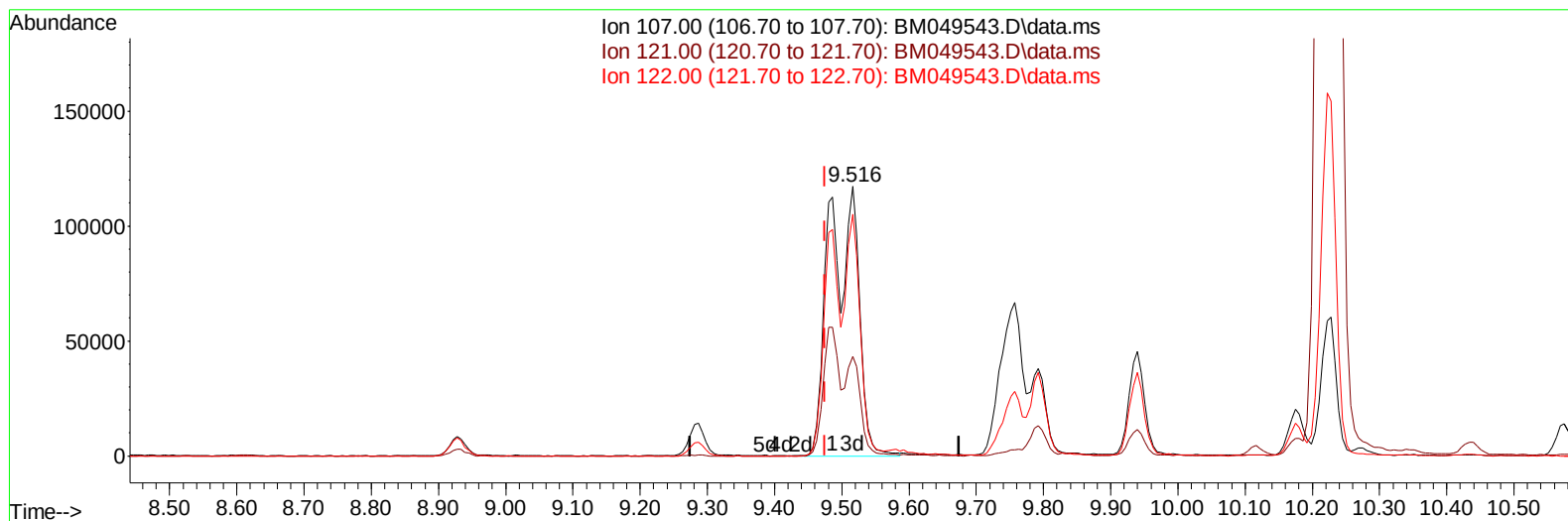
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(26) 2,4-Dimethylphenol

9.516min (+ 0.041) 18.72 ng/ul m

response 356283

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	49.50	36.87#
122.00	85.90	89.65
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Quant Time: Jan 31 09:30:24 2025
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 QLast Update : Tue Jan 28 14:45:18 2025
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.510	152	292405	20.000	ng/ul	0.00
20) Naphthalene-d8	10.269	136	1116302	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.151	164	777423	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.910	188	1608254	20.000	ng/ul	0.00
79) Chrysene-d12	21.139	240	1594680	20.000	ng/ul	0.00
88) Perylene-d12	23.921	264	1726770	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.734	99	1962	0.083	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	6.863	67	4315	0.266	ng/ul	0.00
11) 2-Chlorophenol-d4	7.063	132	4712	0.257	ng/ul	0.01
15) 4-Methylphenol-d8	8.245	113	3167	0.182	ng/ul	0.02
21) Nitrobenzene-d5	8.663	128	2637	0.309	ng/ul	0.00
24) 2-Nitrophenol-d4	9.369	143	2141	0.230	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.922	165	5709	0.300	ng/ul	0.02
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	13.574	166	14271	0.250	ng/ul	0.00
49) Acenaphthylene-d8	13.839	160	17664	0.264	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.157	176	15072	0.301	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.004	188	30169	0.376	ng/ul	0.00
81) Pyrene-d10	19.315	212	26165	0.281	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.733	264	26342	0.286	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	6.757	94	189145	7.595	ng/ul	96
18) 4-Methylphenol	8.304	108	59138	3.139	ng/ul	93
26) 2,4-Dimethylphenol	9.516	107	356283m	18.718	ng/ul	
30) Naphthalene	10.316	128	82382	1.386	ng/ul	99

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

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