

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
 Data File : BP010894.D
 Acq On : 28 Jun 2022 01:01
 Operator : CG/JU
 Sample : N3374-01DL 10X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

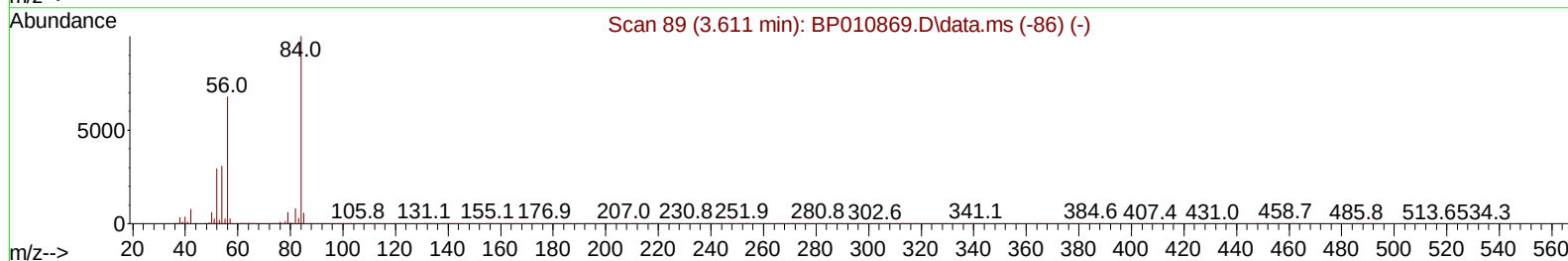
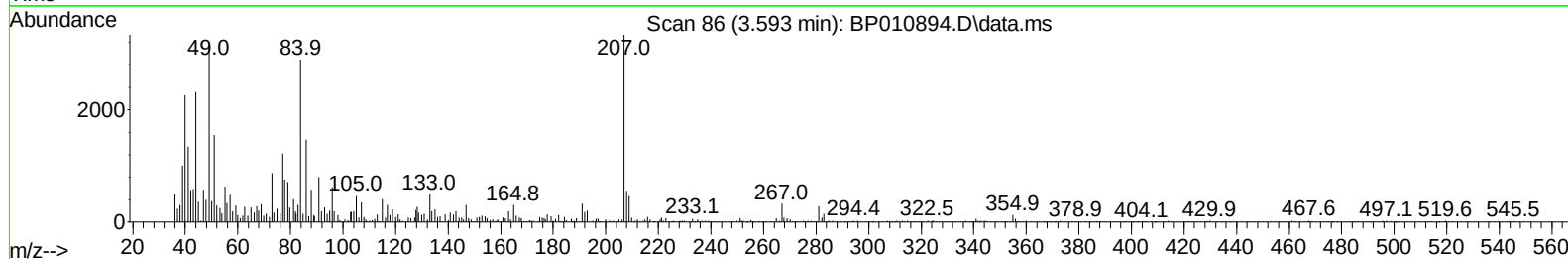
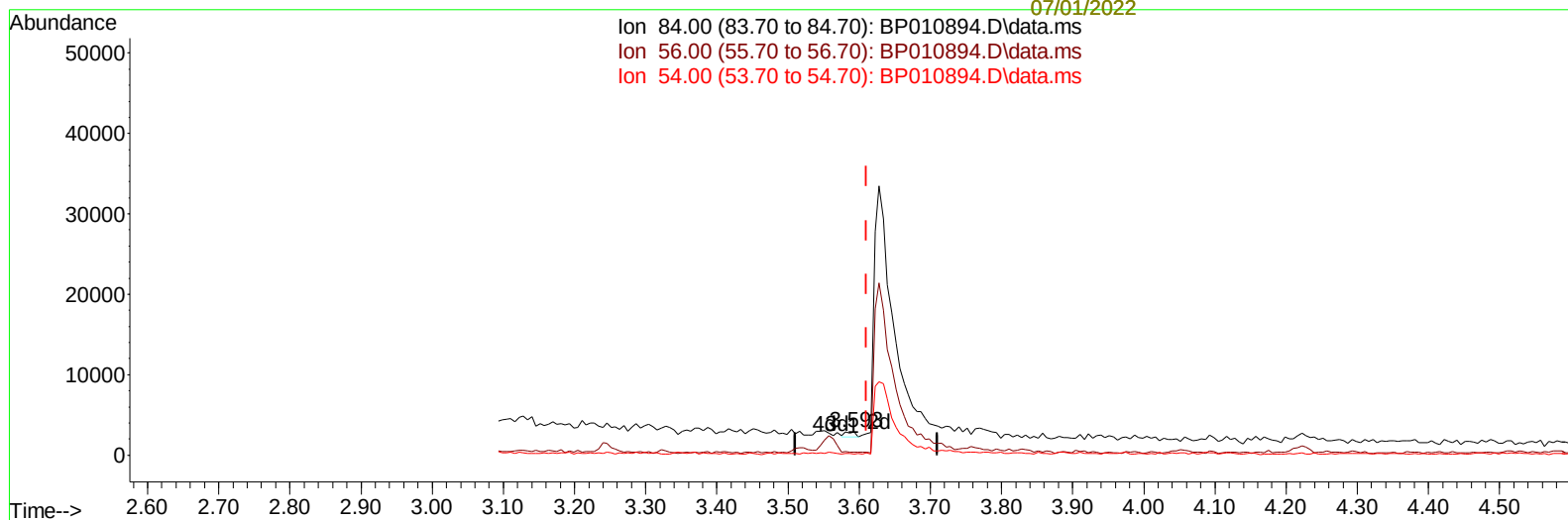
Instrument :
 BNA_P
 ClientSampleId :
 EW4S6DL

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 06/30/2022
 Supervised By :mohammad ahmed 07/01/2022

Quant Time: Jun 29 10:29:21 2022
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 Quant Title : SVOA CALIBRATION
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 ahmed
 07/01/2022



TIC: BP010894.D\data.ms

(4) Pyridine-d5 (S)

3.593min (-0.018) 0.02 ng/ul

response 636

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	11.53#
54.00	44.80	5.63#
0.00	0.00	0.00

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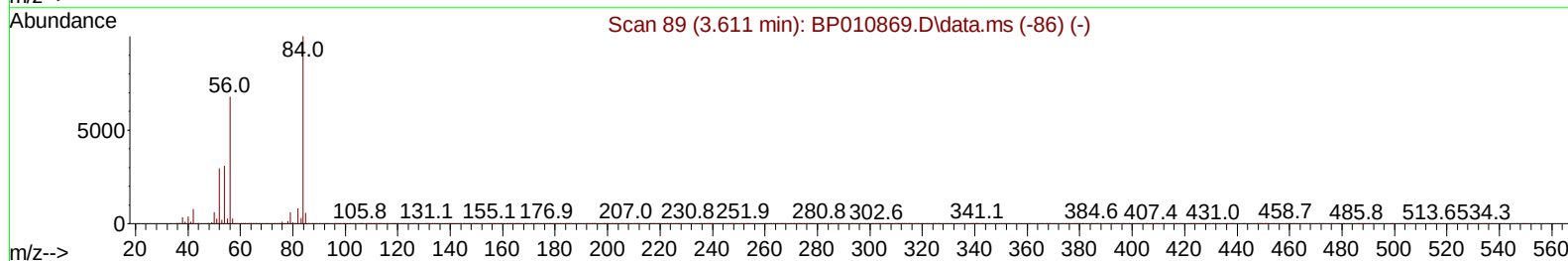
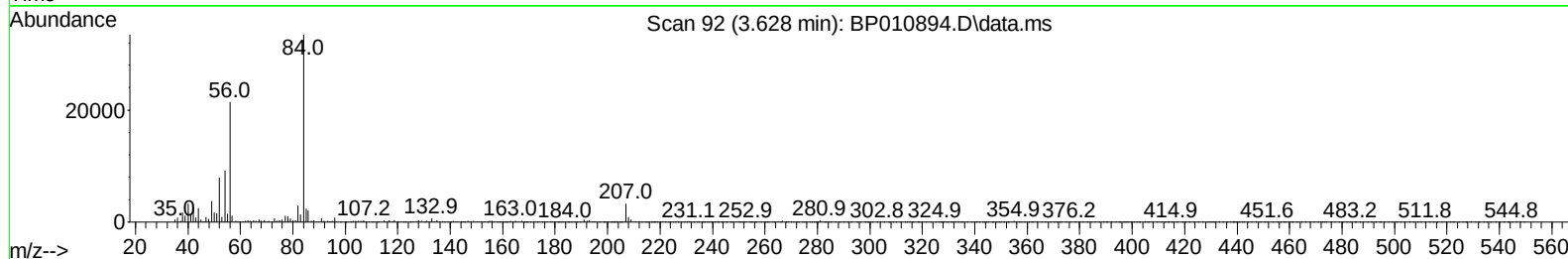
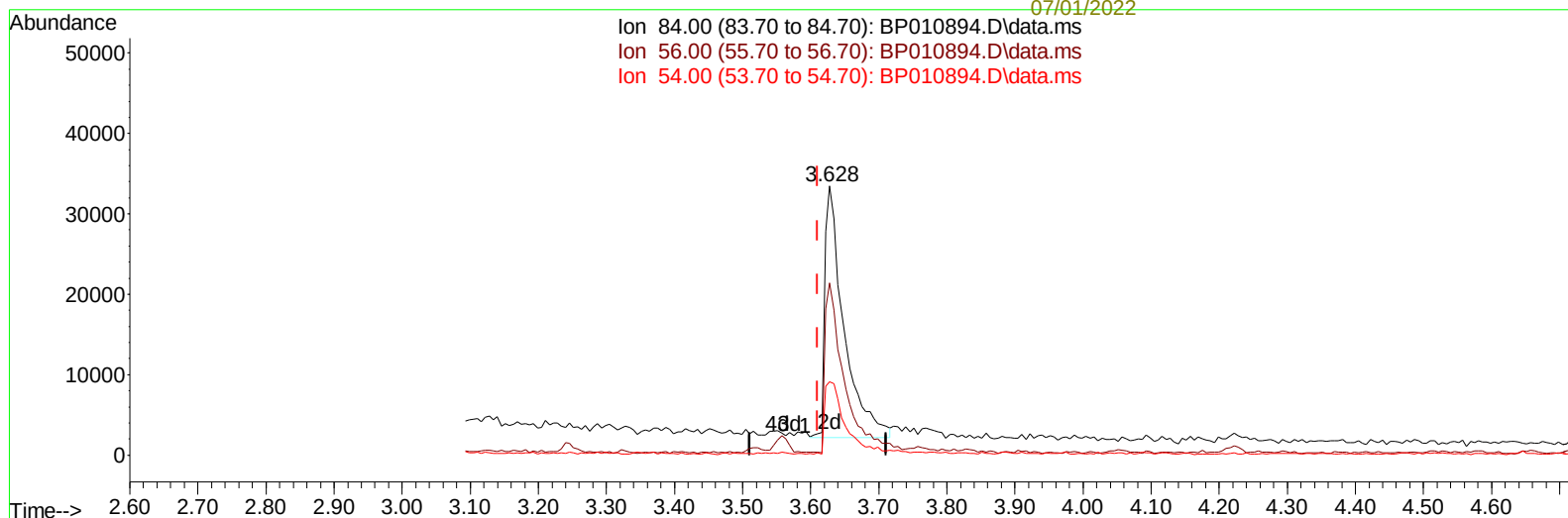
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(4) Pyridine-d5 (S)

3.628min (+ 0.017) 2.03 ng/ul m

response 60689

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	63.98#
54.00	44.80	27.36#
0.00	0.00	0.00

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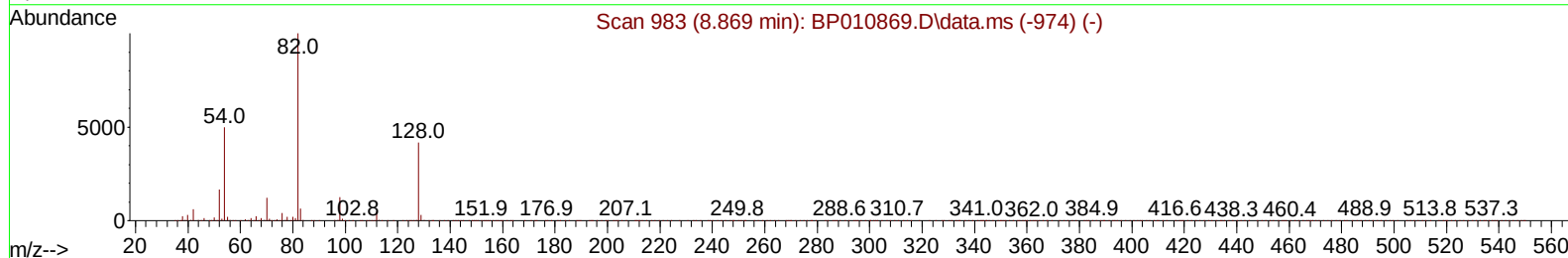
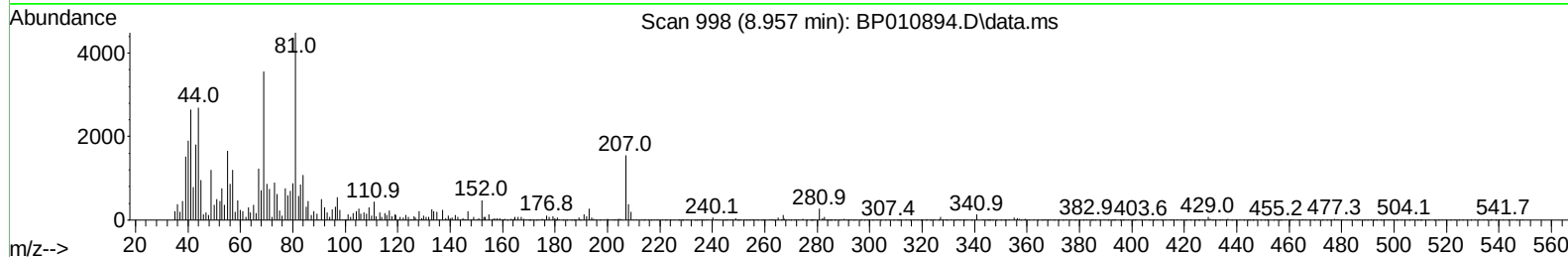
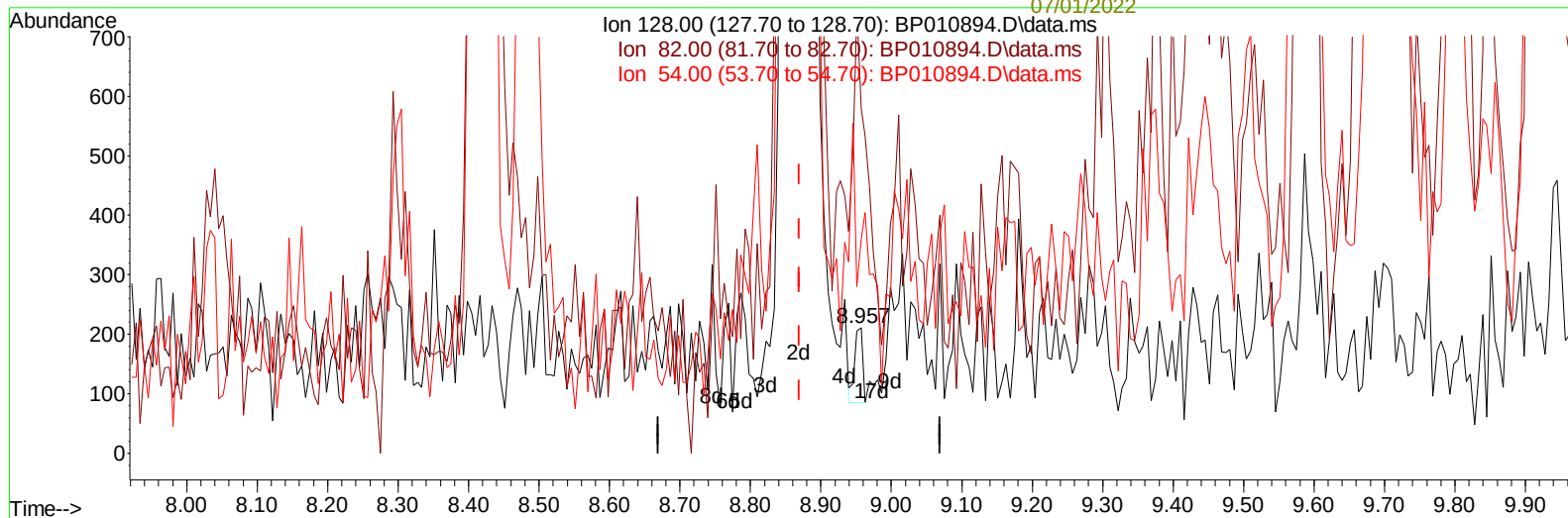
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(21) Nitrobenzene-d5 (S)

8.957min (+ 0.088) 0.01 ng/uL

response 98

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	249.70	274.88
54.00	155.90	170.62
0.00	0.00	0.00

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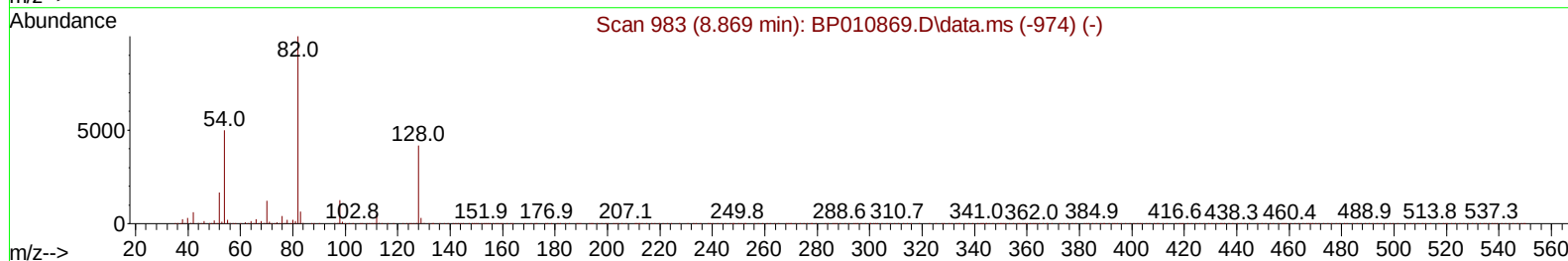
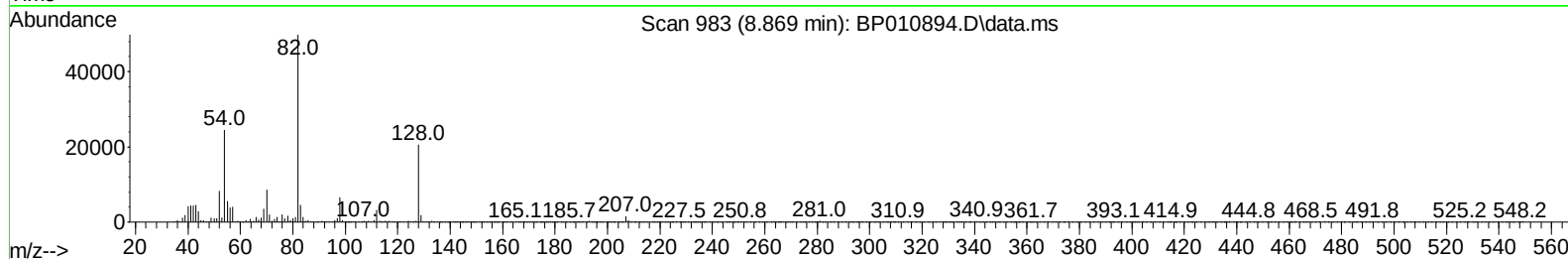
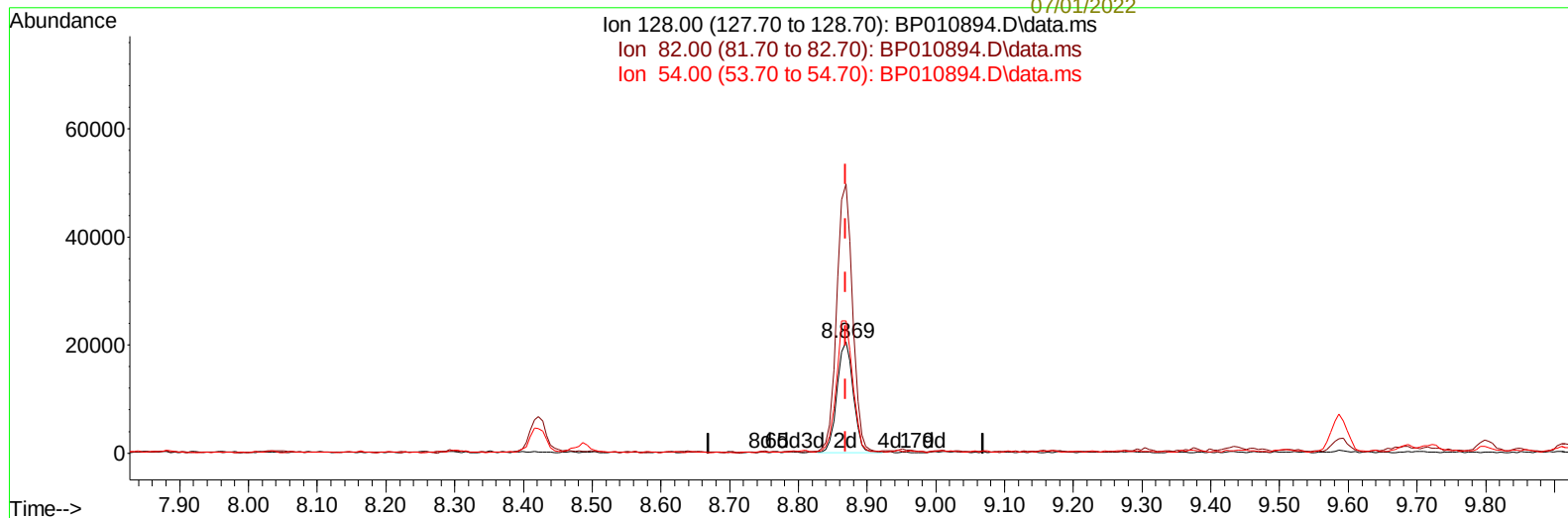
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(21) Nitrobenzene-d5 (S)

8.869min (-0.000) 3.34 ng/ul m

response 33711

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	249.70	241.61
54.00	155.90	118.91#
0.00	0.00	0.00

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Misc :
ALS Vial : 27 Sample Multiplier: 1

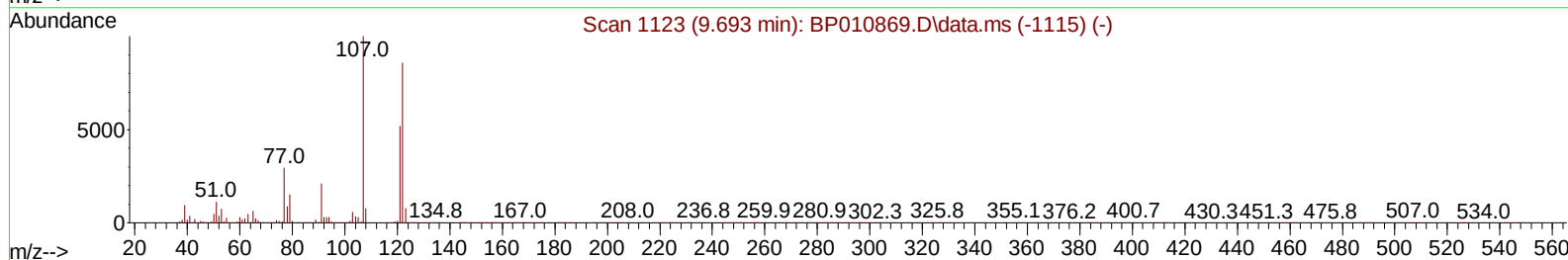
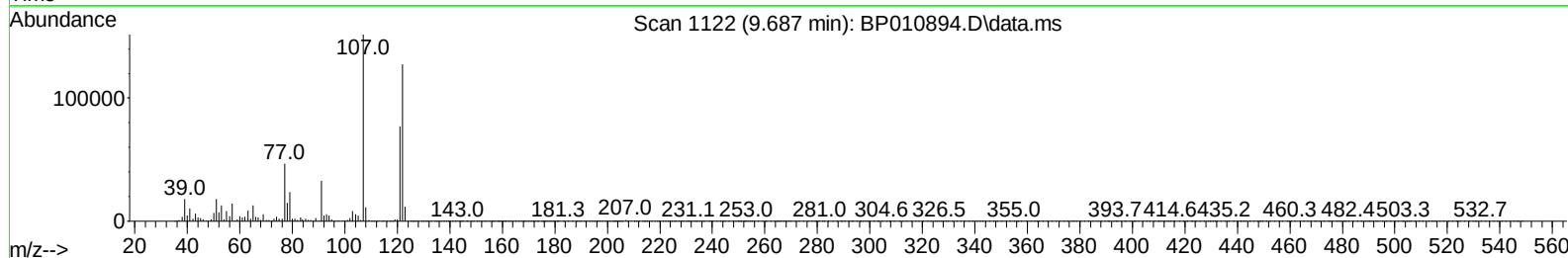
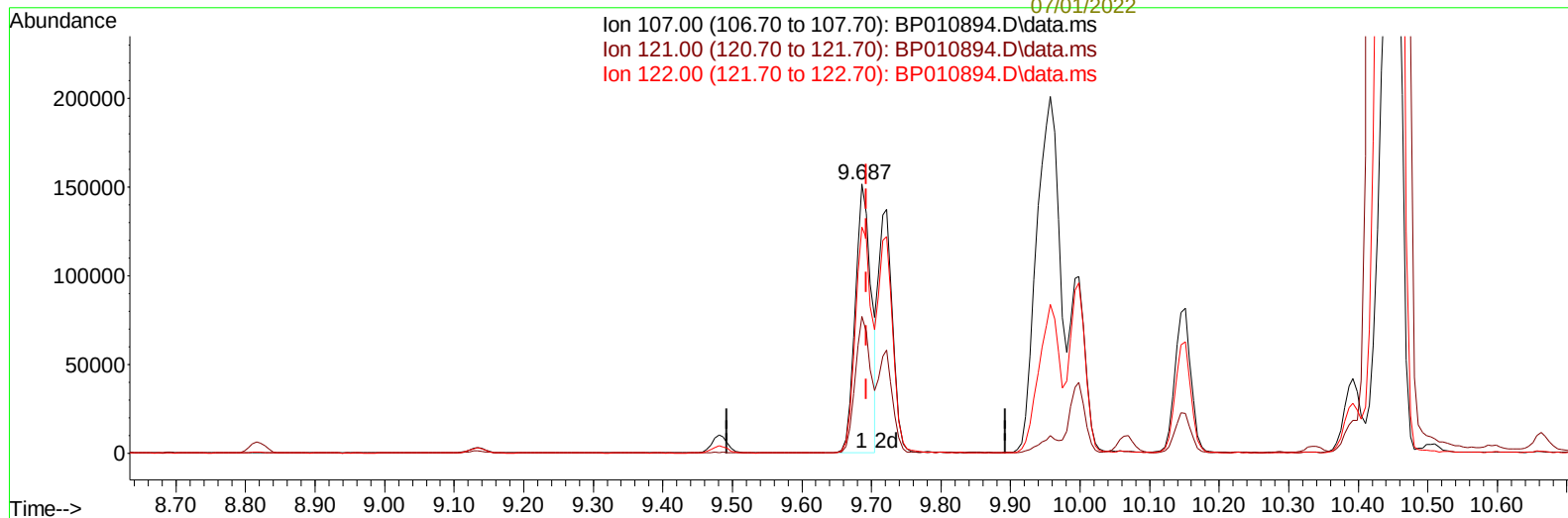
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TIC: BP010894.D\data.ms

(26) 2,4-Dimethylphenol

9.687min (-0.006) 8.63 ng/u1

response 238158

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	48.90	50.77
122.00	82.60	84.16
0.00	0.00	0.00

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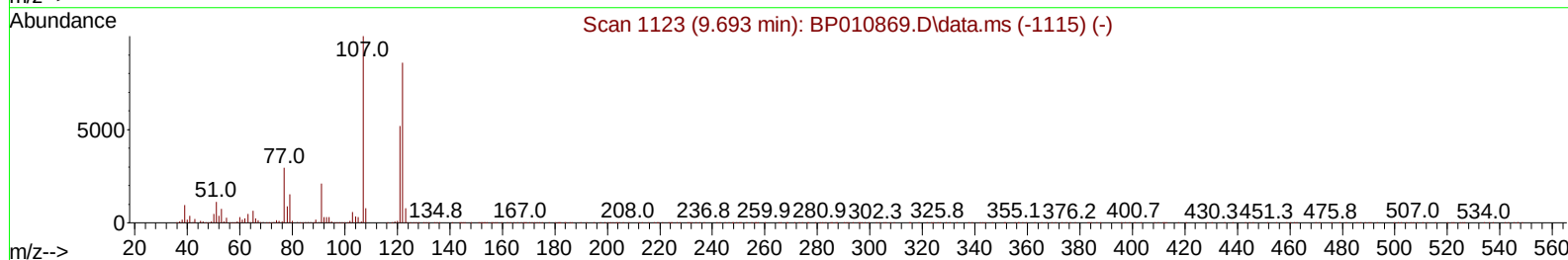
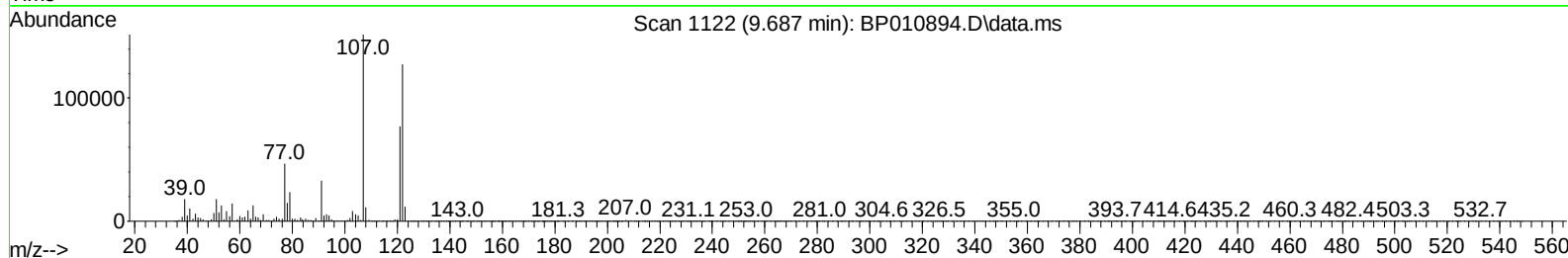
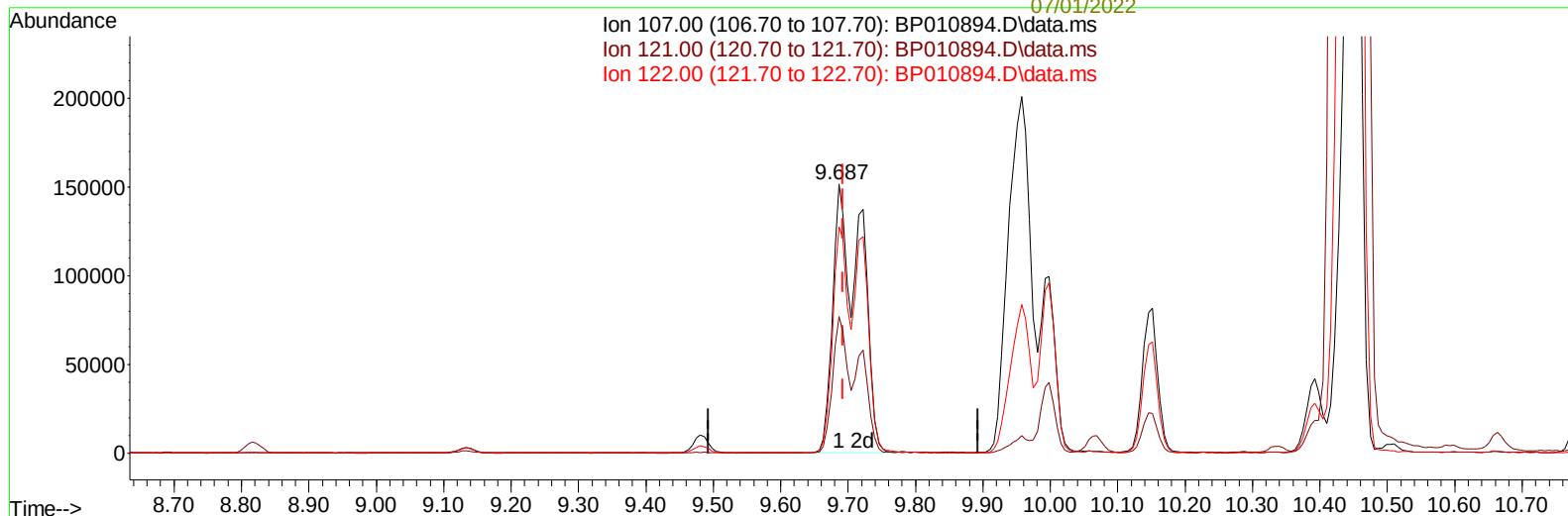
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TIC: BP010894.D\data.ms

(26) 2,4-Dimethylphenol

9.687min (-0.006) 15.66 ng/ul m

response 432017

Ion	Exp%	Act%
107.00	100.00	100.00
121.00	48.90	50.77
122.00	82.60	84.16
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	7.704	152	425102	20.000	ng/ul	0.00
20) Naphthalene-d8	10.504	136	1583631	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.363	164	857124	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.127	188	1602115	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.239	240	1513717	20.000	ng/ul	# 0.00
88) Perylene-d12	23.603	264	1687170	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	3333	0.297	ng/uL	0.00
4) Pyridine-d5	3.628	84	60689m	2.029	ng/ul	0.02
7) Phenol-d5	6.887	99	31410	0.933	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.051	67	64282	2.921	ng/ul	0.00
11) 2-Chlorophenol-d4	7.240	132	88989	3.187	ng/ul	0.00
15) 4-Methylphenol-d8	8.422	113	55552	2.088	ng/ul	0.00
21) Nitrobenzene-d5	8.869	128	33711m	3.344	ng/ul	0.00
24) 2-Nitrophenol-d4	9.586	143	42866	5.139	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.122	165	83718	4.062	ng/ul	0.00
31) 4-Chloroaniline-d4	10.657	131	91192	2.711	ng/ul	0.00
46) Dimethylphthalate-d6	13.786	166	159483	2.716	ng/ul	0.00
49) Acenaphthylene-d8	14.051	160	187729	2.505	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.363	176	148086	2.939	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.480	200	13783	2.597	ng/ul	0.00
73) Anthracene-d10	17.227	188	208834	2.930	ng/ul	0.00
81) Pyrene-d10	19.474	212	234283	2.841	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.445	264	238171	2.857	ng/ul	-0.01
Target Compounds						
					Qvalue	
8) Phenol	6.910	94	1315135	36.540	ng/ul#	85
18) 4-Methylphenol	8.487	108	67473	2.418	ng/ul	96
26) 2,4-Dimethylphenol	9.687	107	432017m	15.656	ng/ul	

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

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