

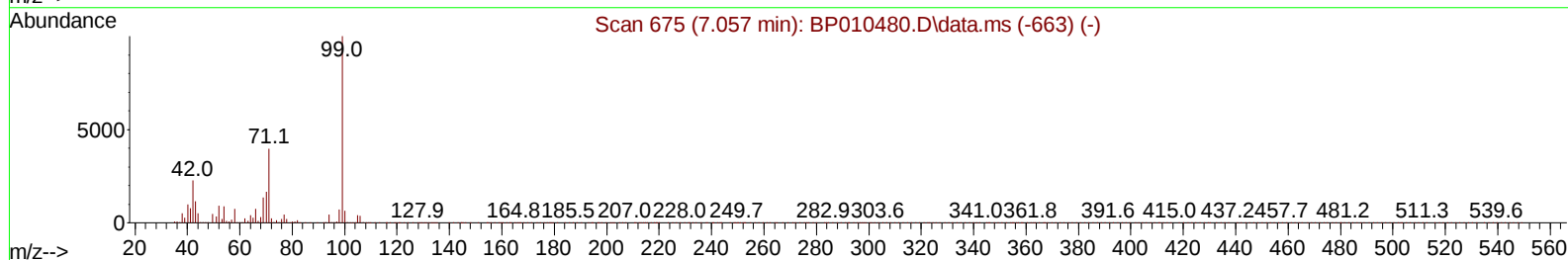
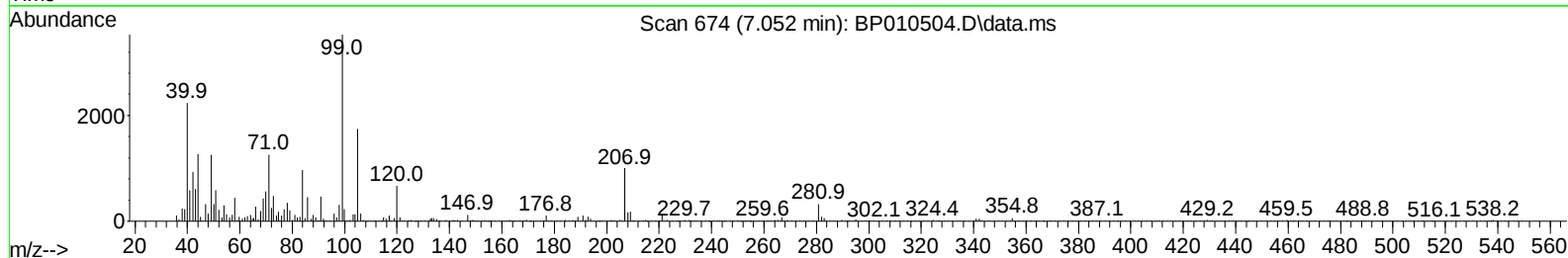
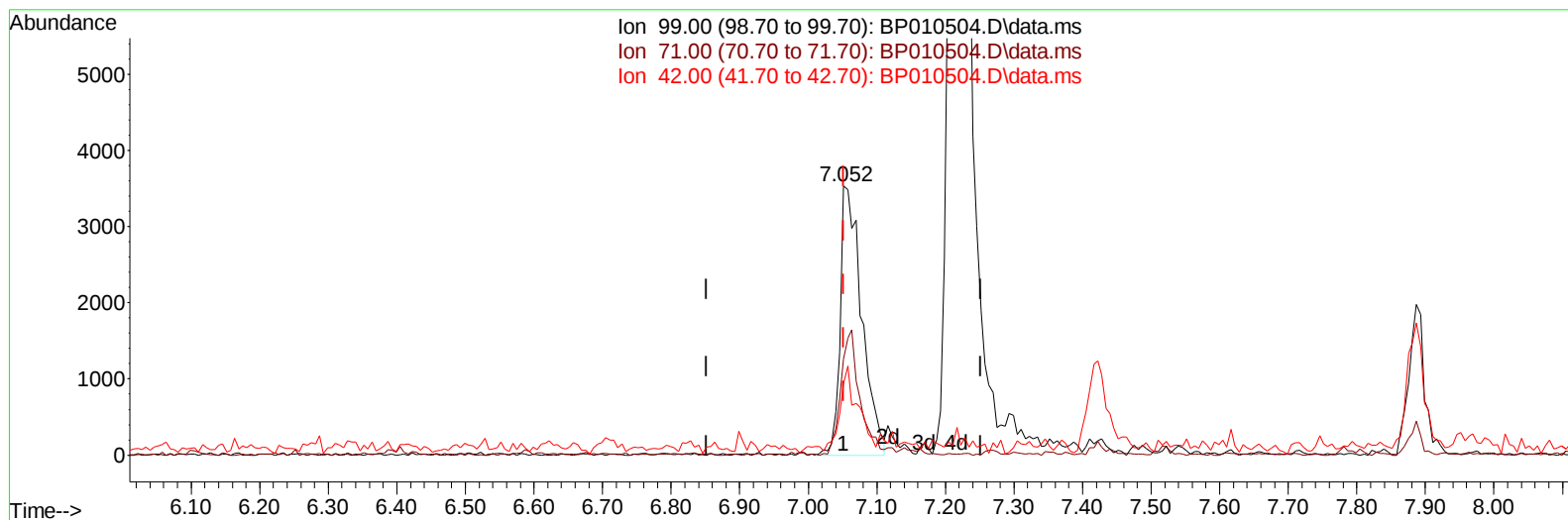
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
 Data File : BP010504.D
 Acq On : 24 May 2022 07:17
 Operator : CG/JU
 Sample : N2918-03DL 10X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBTE0DL

Manual IntegrationsAPPROVED

Quant Time: May 24 07:47:36 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 24 05:50:40 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/24/2022
 Supervised By :Yogesh Patel 05/26/2022



TIC: BP010504.D\data.ms

(7) Phenol-d5 (S)

7.052min (+ 0.000) 0.49 ng/u1

response 7650

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	33.50	35.72
42.00	7.30	26.49#
0.00	0.00	0.00

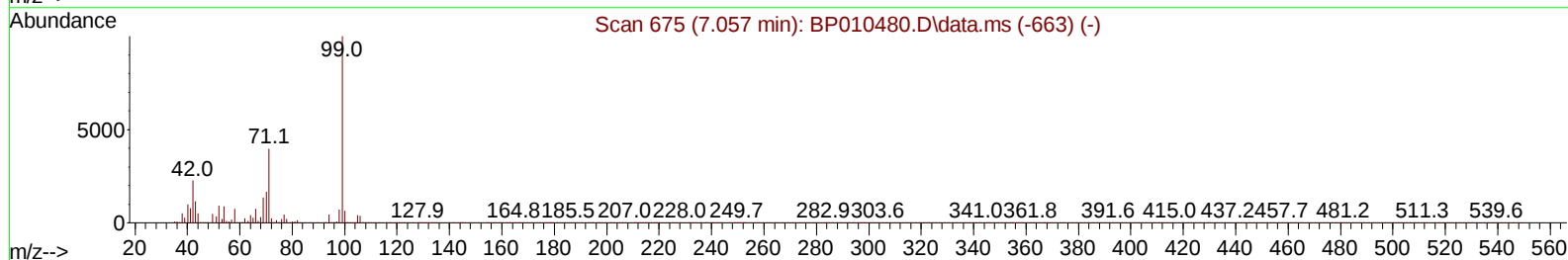
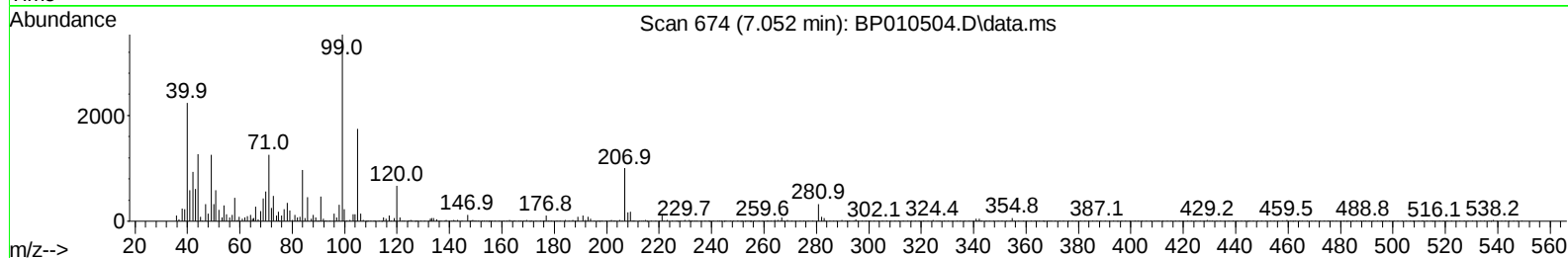
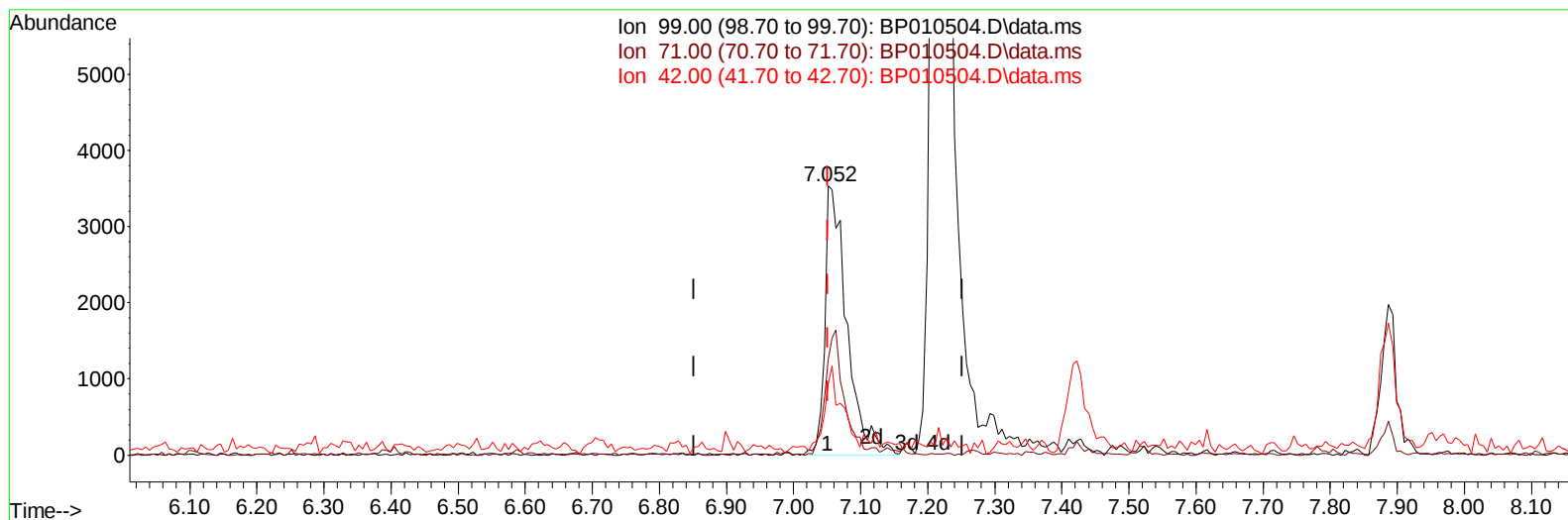
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 Supervised By :Yogesh Patel 05/26/2022



TIC: BP010504.D\data.ms

(7) Phenol-d5 (S)

7.052min (+ 0.000) 0.51 ng/ul m

response 8060

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	33.50	35.72
42.00	7.30	26.49#
0.00	0.00	0.00

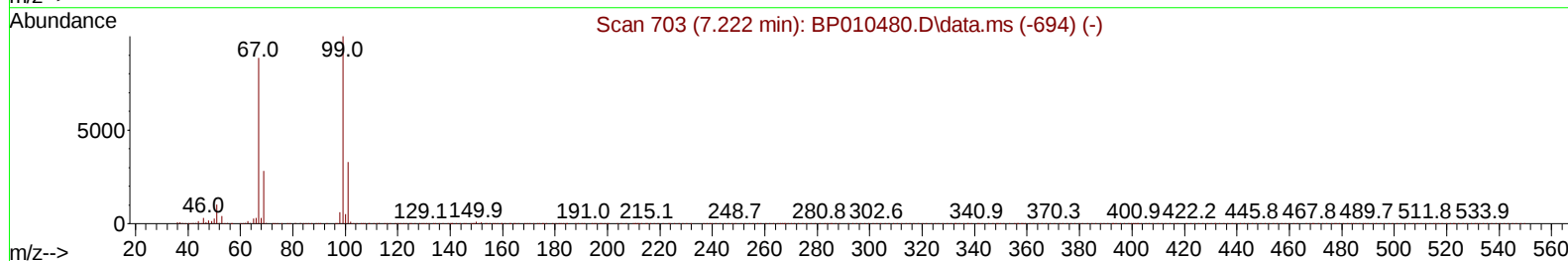
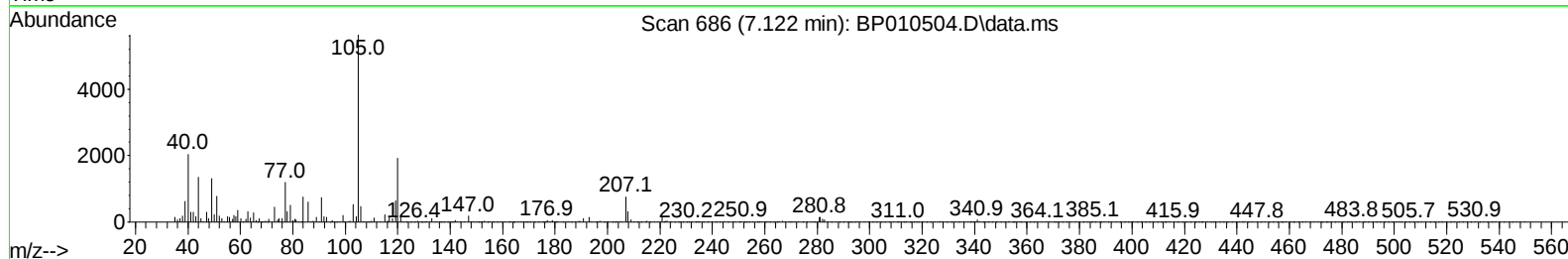
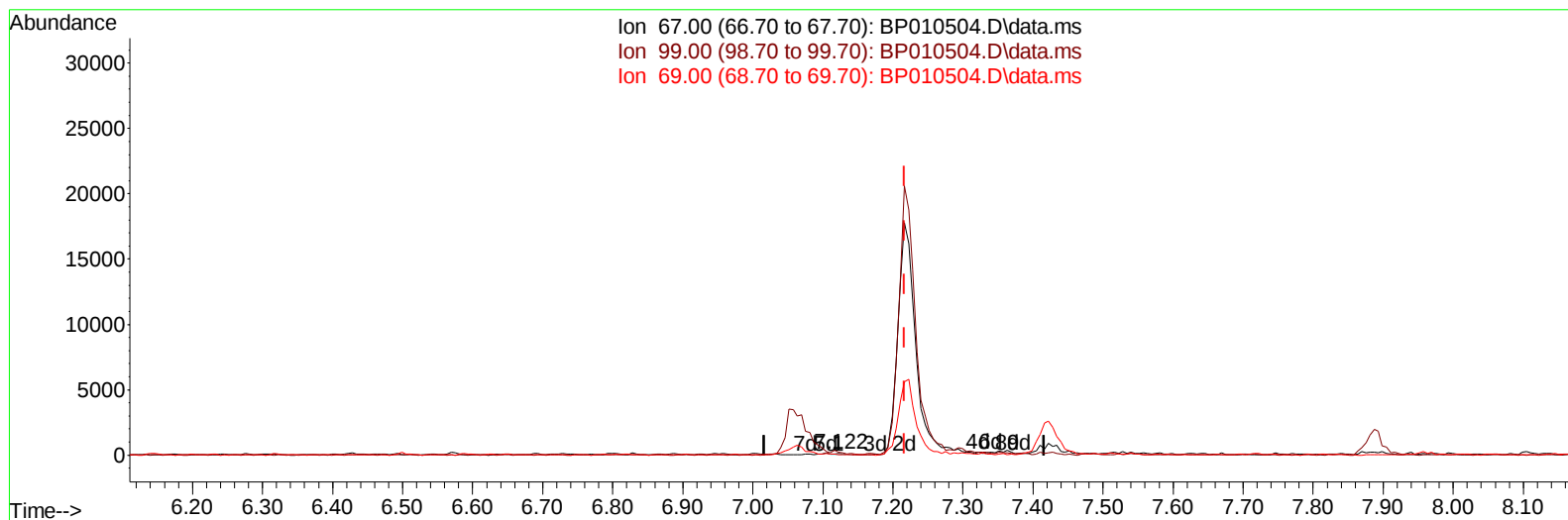
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
 Data File : BP010504.D
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 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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TIC: BP010504.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.122min (-0.094) 0.01 ng/u1

response 92

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	192.62
69.00	40.10	46.72
0.00	0.00	0.00

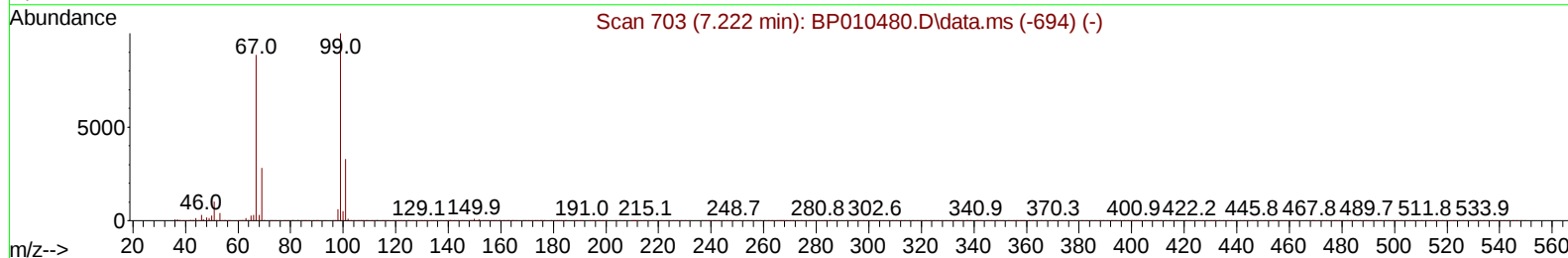
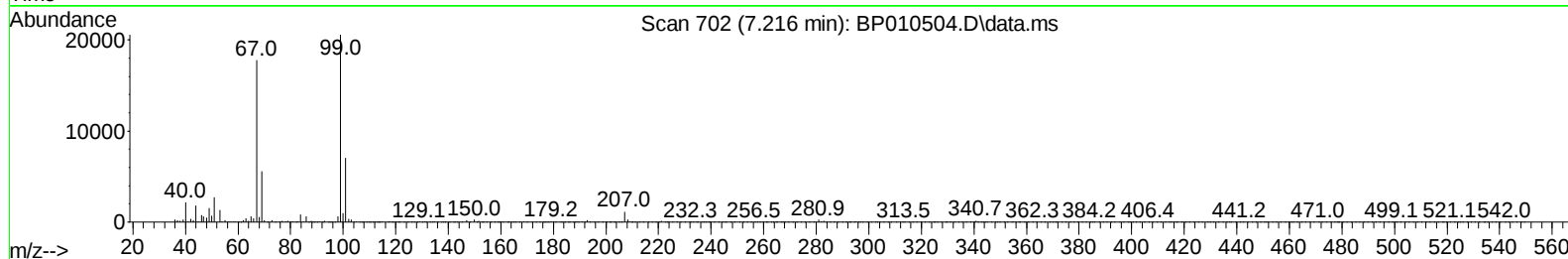
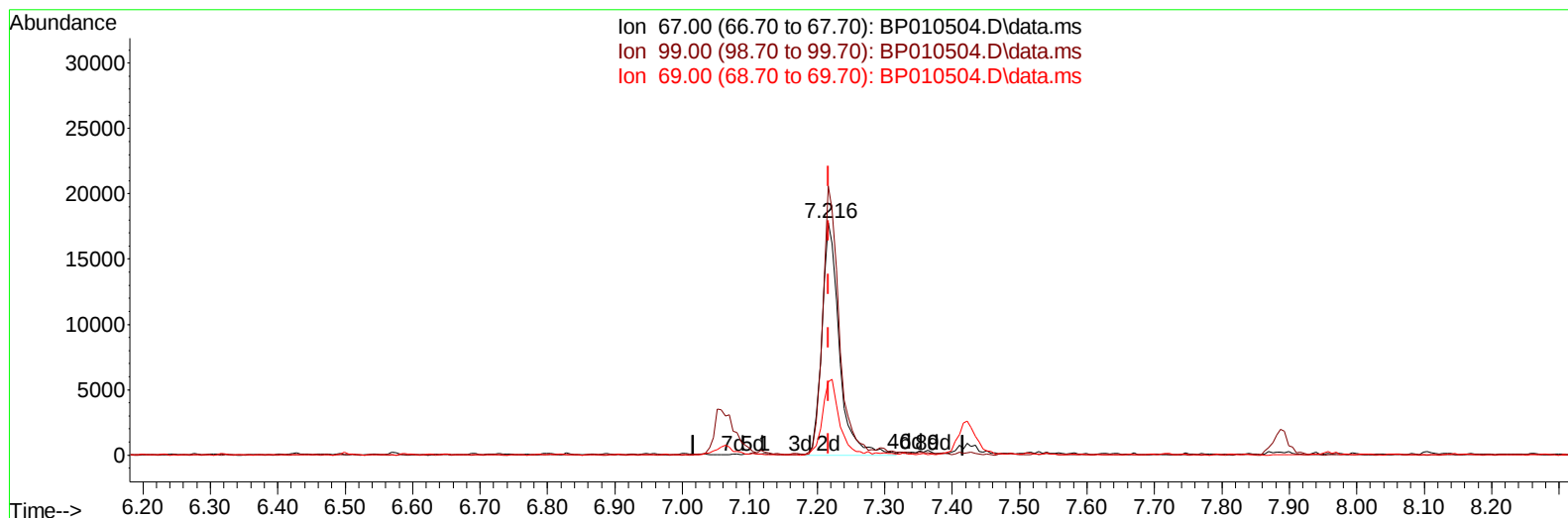
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Instrument :
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TIC: BP010504.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.216min (+ 0.000) 3.01 ng/ul m

response 31799

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	115.80#
69.00	40.10	31.51#
0.00	0.00	0.00

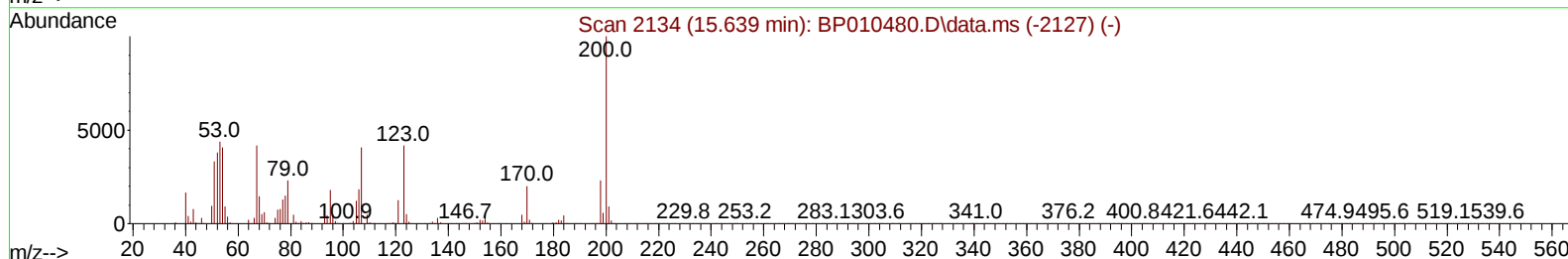
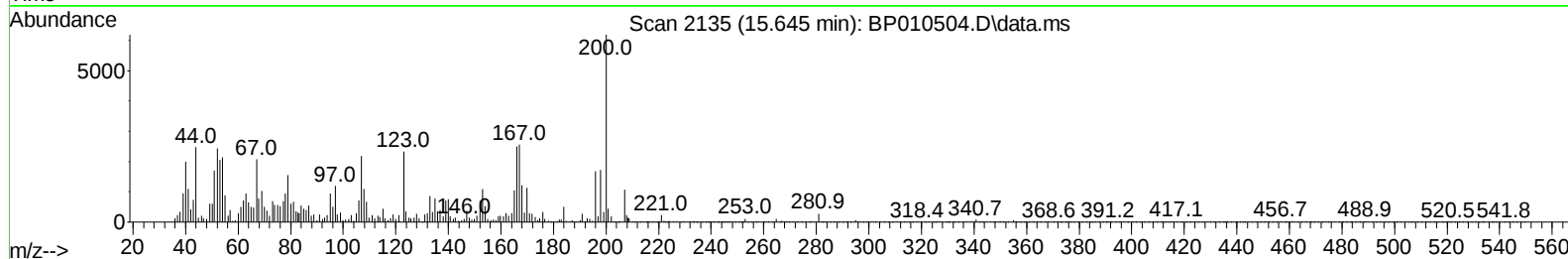
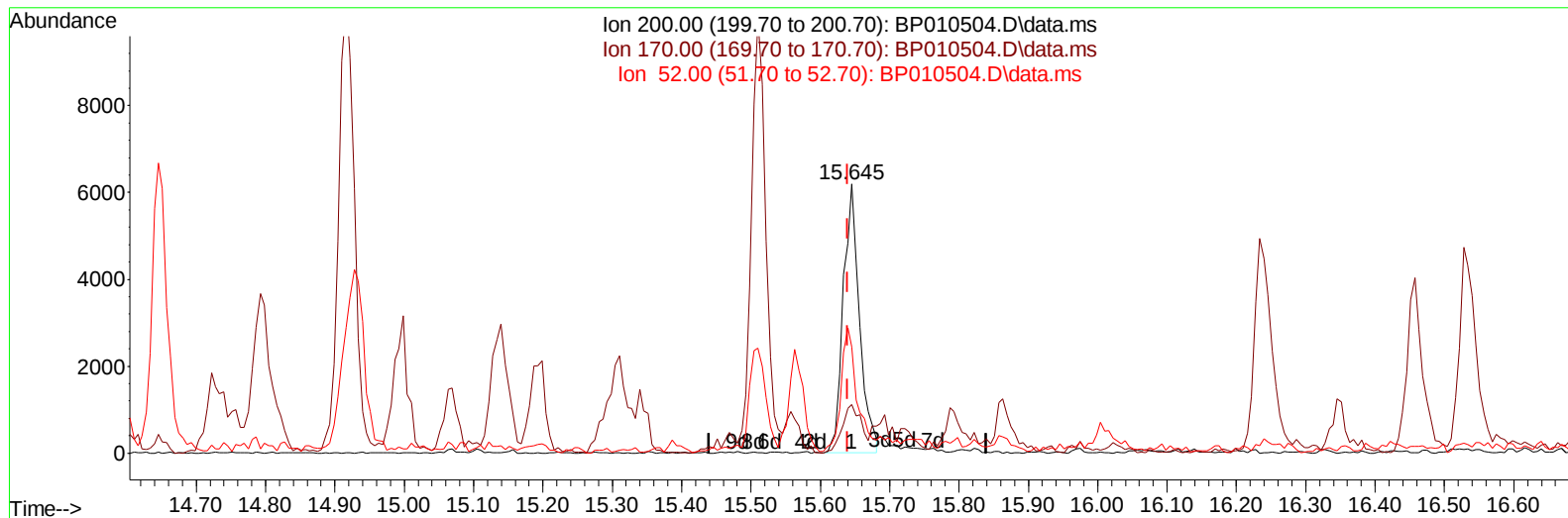
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 ALS Vial : 33 Sample Multiplier: 1

Instrument :
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TIC: BP010504.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.645min (+ 0.006) 1.38 ng/u1

response 9995

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	18.16#
52.00	21.20	39.49#
0.00	0.00	0.00

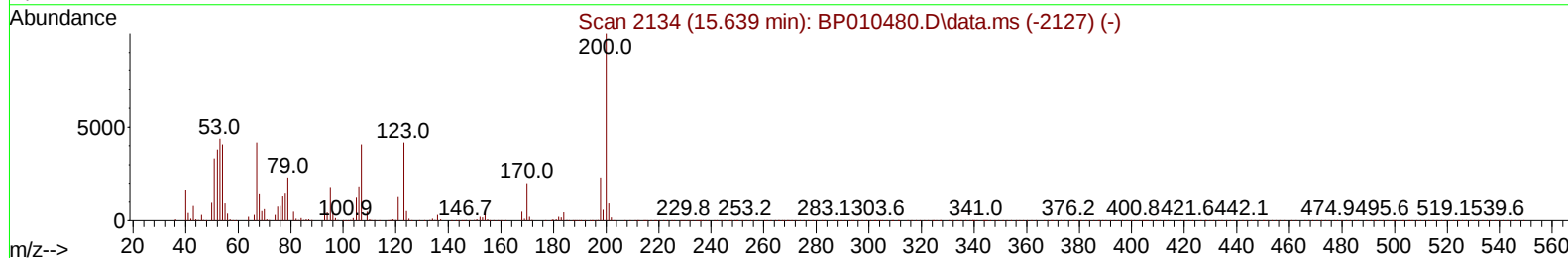
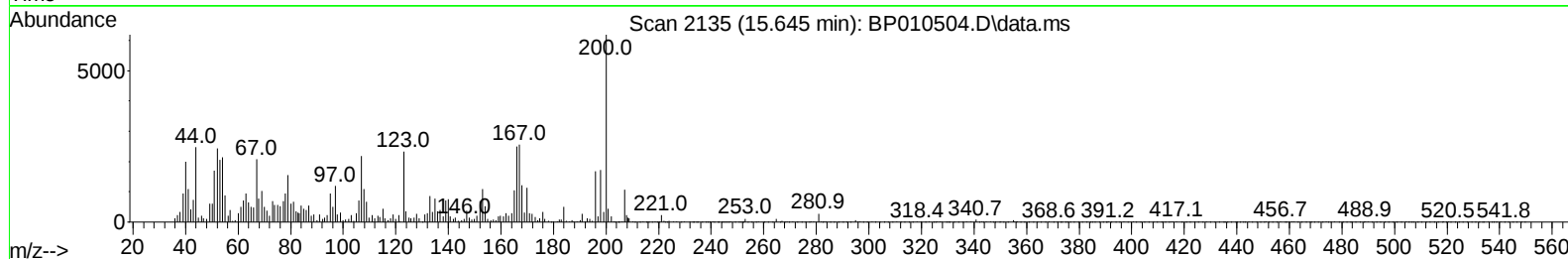
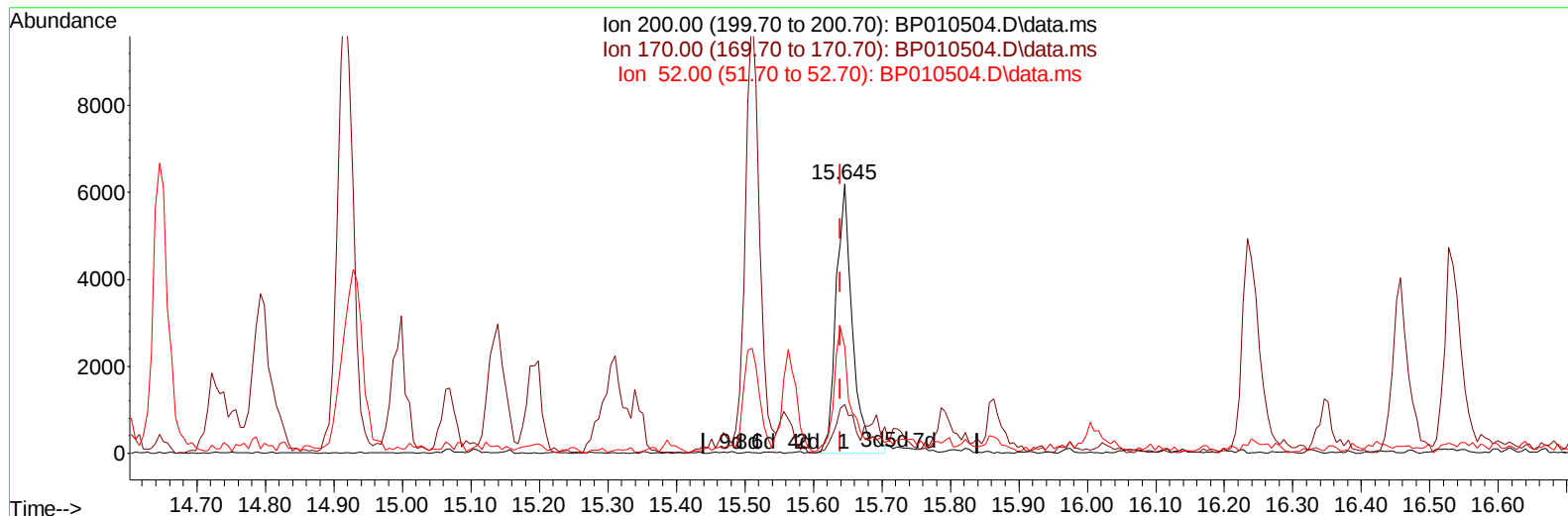
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Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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QLast Update : Tue May 24 05:50:40 2022
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Supervised By :Yogesh Patel 05/26/2022



TIC: BP010504.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.645min (+ 0.006) 1.44 ng/ul m

response 10393

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	18.16#
52.00	21.20	39.49#
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	7.887	152	187156	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	844656	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.516	164	573044	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.269	188	1234776	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.351	240	1071804	20.000	ng/ul	# 0.00
88) Perylene-d12	23.774	264	822160	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	1180	0.269	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.052	99	8060m	0.513	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	31799m	3.013	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132	24024	1.999	ng/ul	0.00
15) 4-Methylphenol-d8	8.605	113	15835	1.186	ng/ul	0.00
21) Nitrobenzene-d5	9.052	128	18590	2.844	ng/ul	0.00
24) 2-Nitrophenol-d4	9.775	143	17364	2.506	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	30139	2.301	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	49048	2.518	ng/ul	0.00
46) Dimethylphthalate-d6	13.928	166	130245	3.085	ng/ul	0.00
49) Acenaphthylene-d8	14.210	160	145734	2.993	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.510	176	116206	3.164	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.369	188	181671	3.249	ng/ul	0.00
81) Pyrene-d10	19.598	212	216196	3.798	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.616	264	128078	3.102	ng/ul	0.00
Target Compounds						
					Qvalue	
30) Naphthalene	10.752	128	19630352	444.954	ng/ul#	95
36) 2-Methylnaphthalene	12.346	142	1844319	59.051	ng/ul	90
37) 1-Methylnaphthalene	12.563	142	1003758	31.830	ng/ul#	94
43) 1,1'-Biphenyl	13.351	154	430440	10.234	ng/ul#	91
50) Acenaphthylene	14.240	152	92256	1.751	ng/ul	95
52) Acenaphthene	14.581	153	1440468	41.913	ng/ul	98
56) Dibenzofuran	14.916	168	1042921	20.707	ng/ul	98
61) Fluorene	15.569	166	685454	16.642	ng/ul	92
71) Pentachlorophenol	16.928	266	23613	2.573	ng/ul#	83
72) Phenanthrene	17.310	178	912763	13.851	ng/ul	99
77) Carbazole	17.675	167	2400904	41.325	ng/ul#	95
80) Fluoranthene	19.275	202	74892	1.132	ng/ul#	95

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

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