

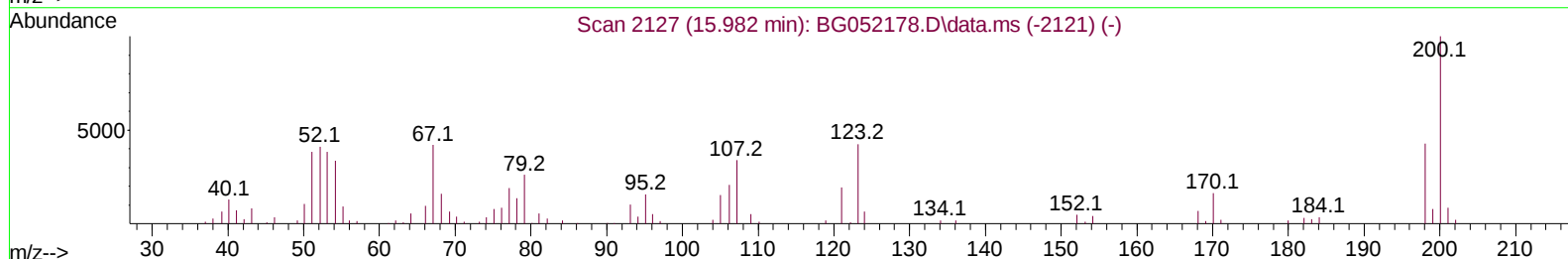
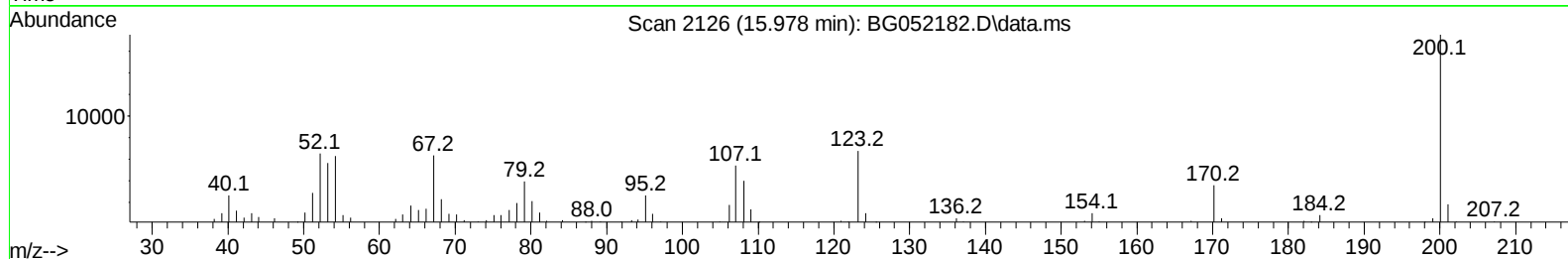
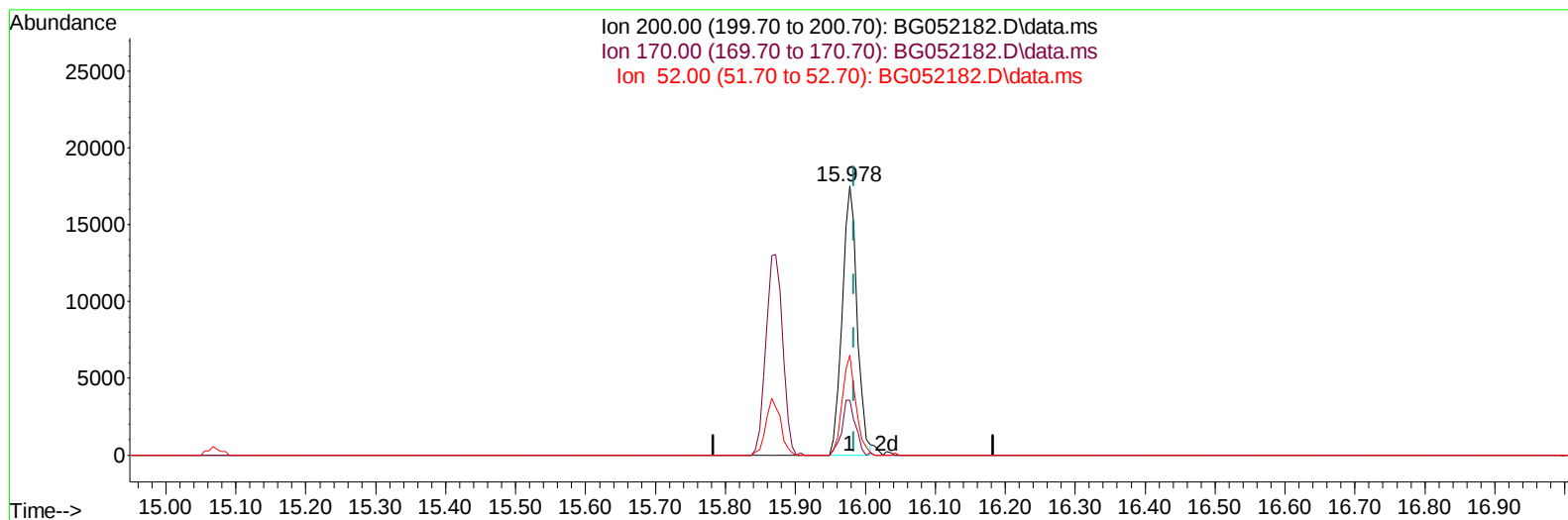
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052182.D
 Acq On : 27 Jan 2022 12:57
 Operator : CG/JU
 Sample : N1279-02
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0VD9

Manual IntegrationsAPPROVED

Quant Time: Jan 29 00:42:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 26 23:58:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/31/2022
 Supervised By :mohammad ahmed 01/31/2022



TIC: BG052182.D\data.ms

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

15.978min (-0.006) 21.91 ng/ul

response 26154

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	20.45
52.00	47.40	37.19#
0.00	0.00	0.00

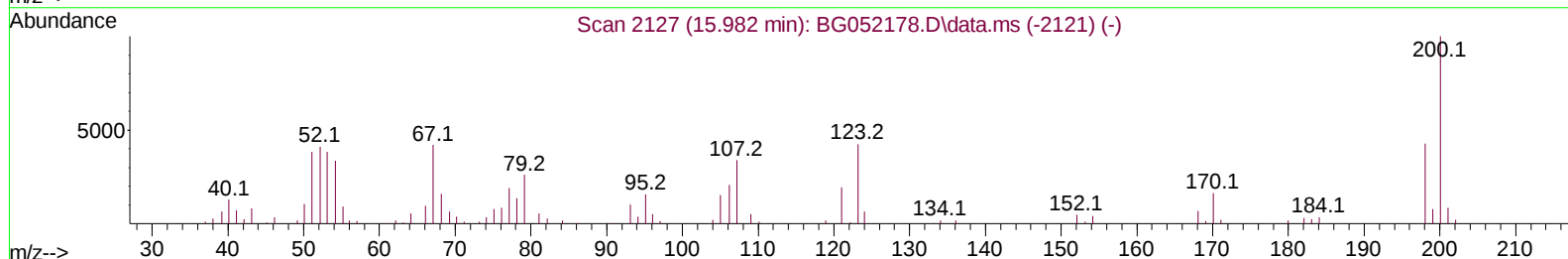
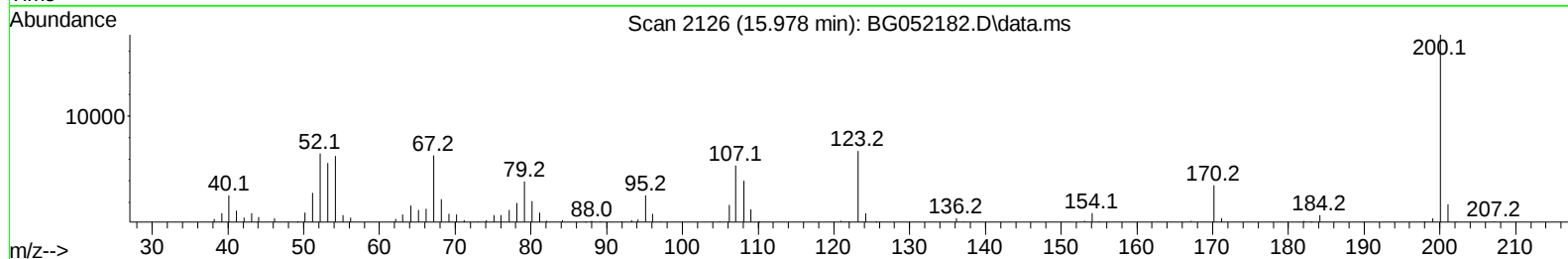
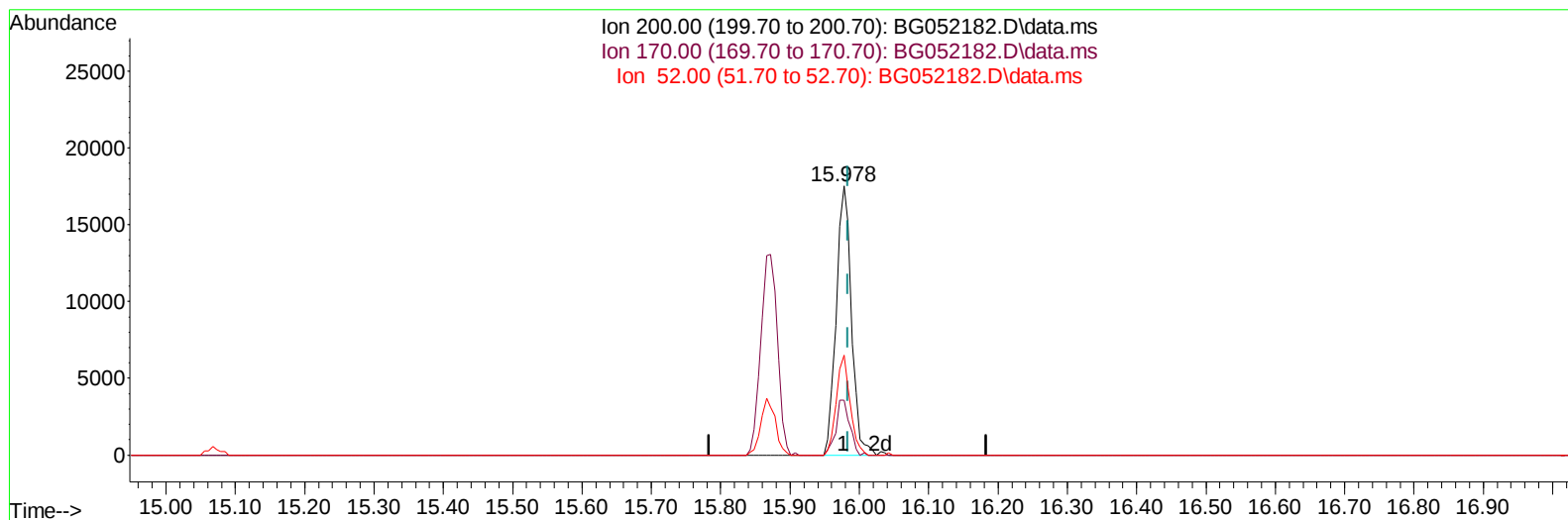
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052182.D
 Acq On : 27 Jan 2022 12:57
 Operator : CG/JU
 Sample : N1279-02
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COVID9

Manual IntegrationsAPPROVED

Quant Time: Jan 27 23:58:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 26 23:58:52 2022
 Response via : Initial Calibration

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TIC: BG052182.D\data.ms

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

15.978min (-0.006) 22.18 ng/ul m

response 26469

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	20.45
52.00	47.40	37.19#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052182.D
 Acq On : 27 Jan 2022 12:57
 Operator : CG/JU
 Sample : N1279-02
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COVID9

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/31/2022
 Supervised By :mohammad ahmed 01/31/2022

Quant Time: Jan 27 23:58:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 26 23:58:52 2022
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.235	152	23194	20.000	ng/ul	0.00
20) Naphthalene-d8	11.073	136	98997	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.879	164	73279	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.628	188	189184	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.946	240	195290	20.000	ng/ul	# 0.00
88) Perylene-d12	25.424	264	194645	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.565	96	2492	3.839	ng/uL	0.00
4) Pyridine-d5	4.000	84	7843	3.887	ng/ul	0.00
7) Phenol-d5	7.383	99	14915	6.570	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	43096	29.906	ng/ul	0.00
11) 2-Chlorophenol-d4	7.759	132	36467	24.263	ng/ul	-0.01
15) 4-Methylphenol-d8	8.934	113	26674	15.116	ng/ul	-0.01
21) Nitrobenzene-d5	9.404	128	26939	34.438	ng/ul	0.00
24) 2-Nitrophenol-d4	10.139	143	28914	33.421	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.691	165	50434	29.824	ng/ul	0.00
31) 4-Chloroaniline-d4	11.196	131	59315	25.187	ng/ul	-0.01
46) Dimethylphthalate-d6	14.268	166	201025	33.188	ng/ul	0.00
49) Acenaphthylene-d8	14.574	160	249897	34.155	ng/ul	0.00
54) 4-Nitrophenol-d4	15.067	143	4865	4.804	ng/ul	0.00
60) Fluorene-d10	15.872	176	182502	34.558	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.978	200	26469m	22.178	ng/ul	0.00
73) Anthracene-d10	17.728	188	316484	35.265	ng/ul	0.00
81) Pyrene-d10	20.008	212	385590	34.307	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.183	264	358892	35.024	ng/ul	0.00

Target Compounds Qvalue

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

