

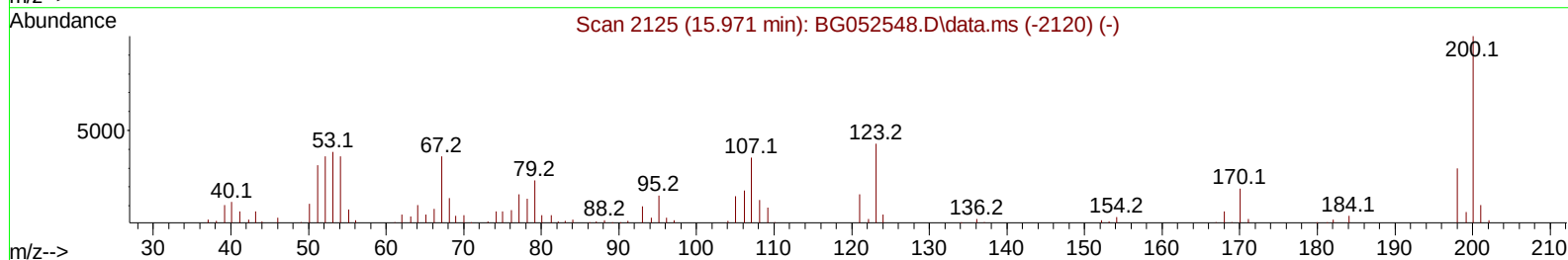
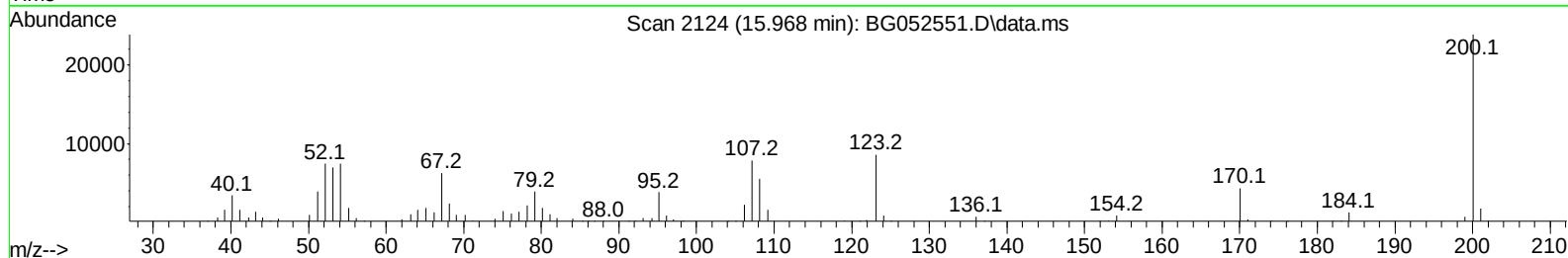
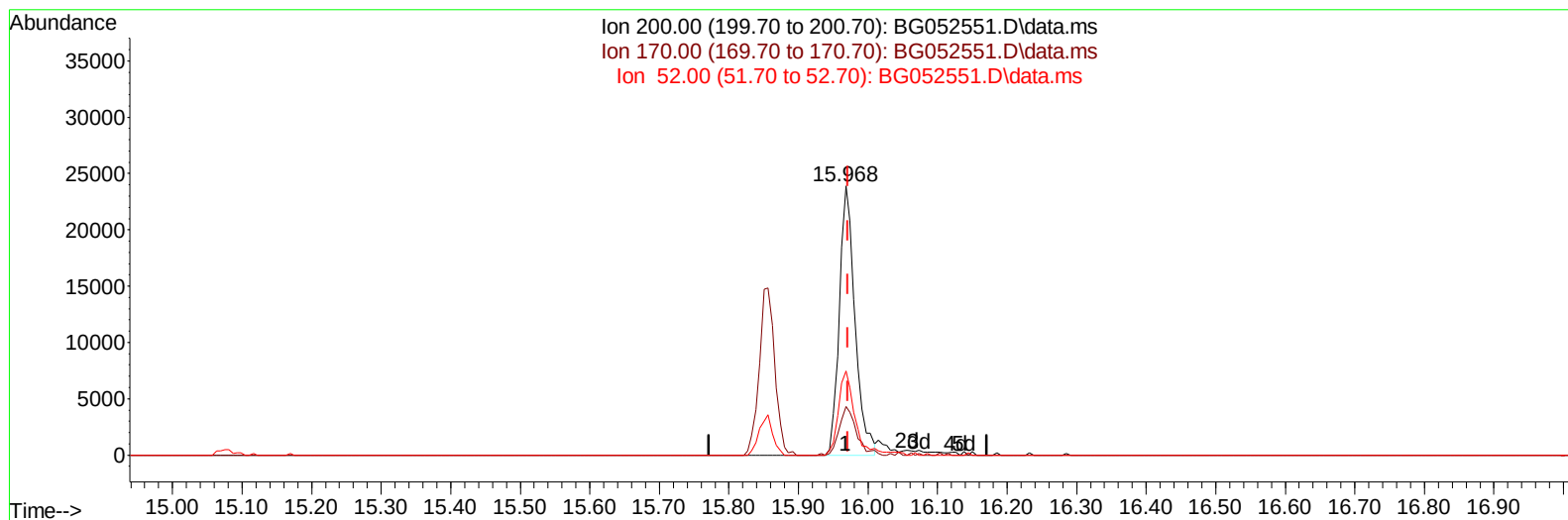
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030322\
Data File : BG052551.D
Acq On : 3 Mar 2022 12:52
Operator : CG/JU
Sample : N1745-05
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
H0AA5

Manual IntegrationsAPPROVED

Quant Time: Mar 04 00:00:49 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022222.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 22 17:32:12 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/04/2022
Supervised By :Yogesh Patel 03/07/2022



TIC: BG052551.D\data.ms

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

15.968min (-0.004) 30.44 ng/u1

response 37577

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	18.17
52.00	47.40	31.21#
0.00	0.00	0.00

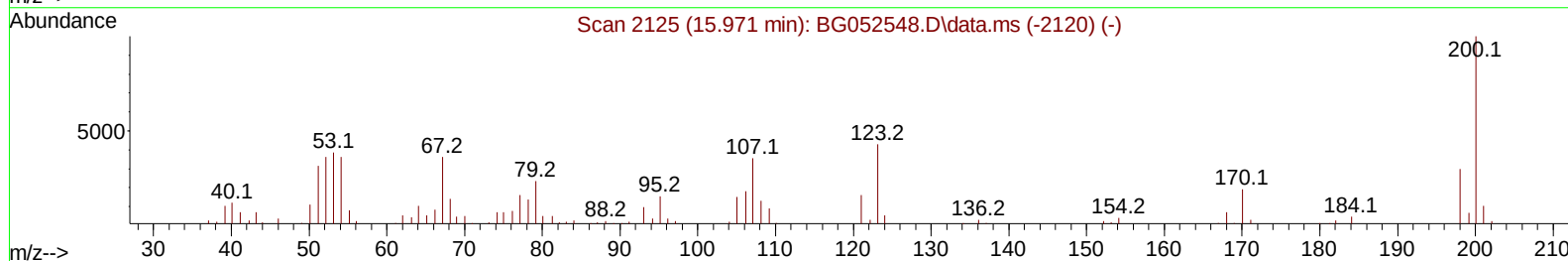
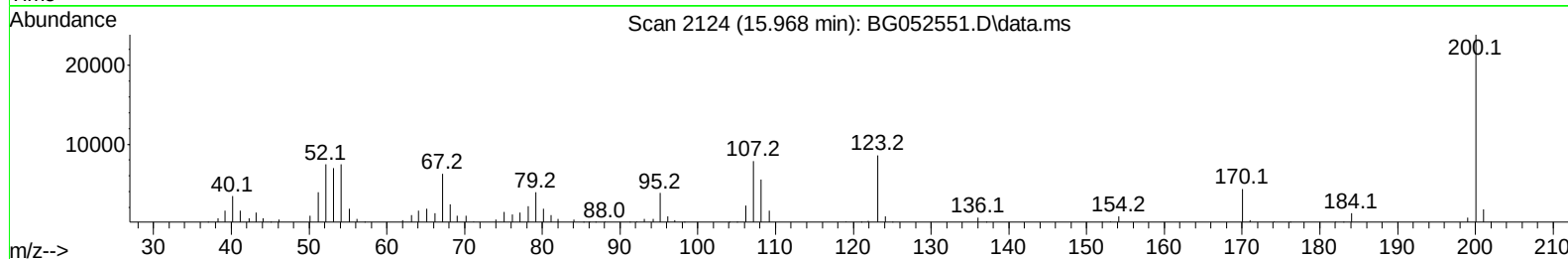
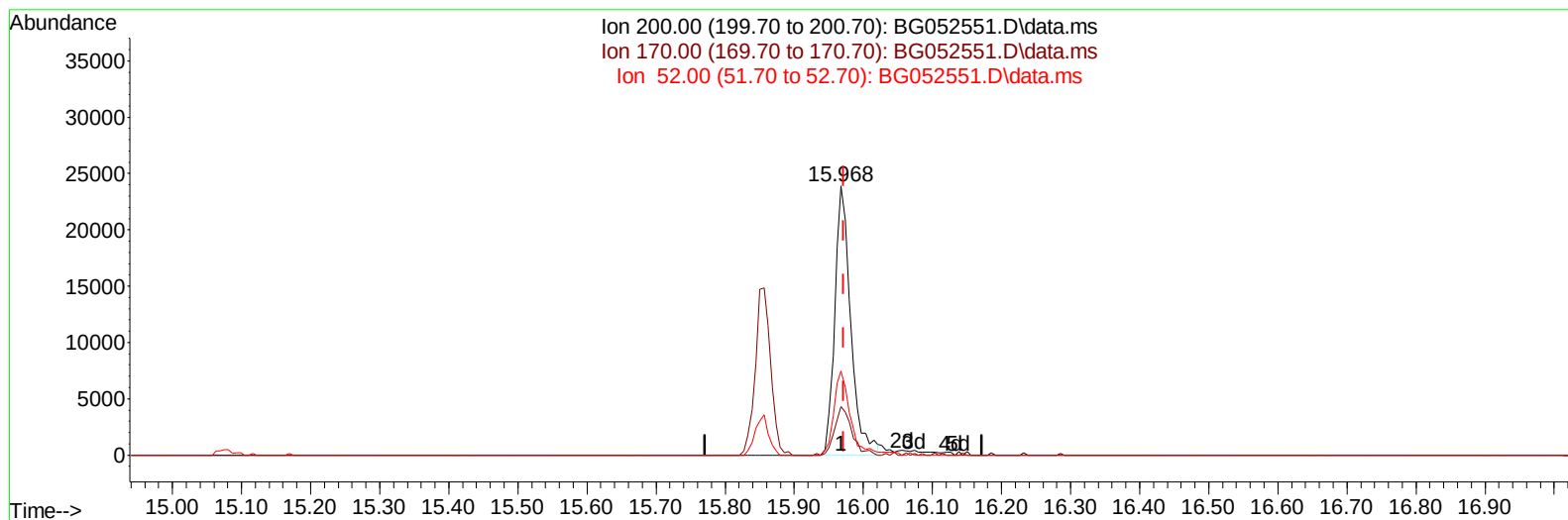
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TIC: BG052551.D\data.ms

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

15.968min (-0.004) 31.08 ng/ul m

response 38376

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	18.17
52.00	47.40	31.21#
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	8.220	152	23335	20.000	ng/ul	0.00
20) Naphthalene-d8	11.057	136	100342	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.864	164	75361	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.613	188	194716	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.930	240	199676	20.000	ng/ul	# 0.00
88) Perylene-d12	25.390	264	200481	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.544	96	2830	4.757	ng/uL	0.00
4) Pyridine-d5	3.984	84	16736	9.326	ng/ul	0.00
7) Phenol-d5	7.368	99	18100	8.180	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.526	67	42174	32.614	ng/ul	0.00
11) 2-Chlorophenol-d4	7.744	132	38891	25.024	ng/ul	0.00
15) 4-Methylphenol-d8	8.925	113	31184	18.113	ng/ul	0.00
21) Nitrobenzene-d5	9.389	128	27489	32.135	ng/ul	0.00
24) 2-Nitrophenol-d4	10.123	143	30507	32.559	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.675	165	52181	29.864	ng/ul	0.00
31) 4-Chloroaniline-d4	11.186	131	68514	28.296	ng/ul	0.00
46) Dimethylphthalate-d6	14.253	166	233883	38.177	ng/ul	0.00
49) Acenaphthylene-d8	14.558	160	264467	34.969	ng/ul	0.00
54) 4-Nitrophenol-d4	15.069	143	7285	7.480	ng/ul	0.01
60) Fluorene-d10	15.856	176	202838	37.689	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.968	200	38376m	31.083	ng/ul	0.00
73) Anthracene-d10	17.713	188	372850	39.636	ng/ul	0.00
81) Pyrene-d10	19.992	212	452192	38.702	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.155	264	432321	40.195	ng/ul	0.00

Target Compounds Qvalue

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

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