

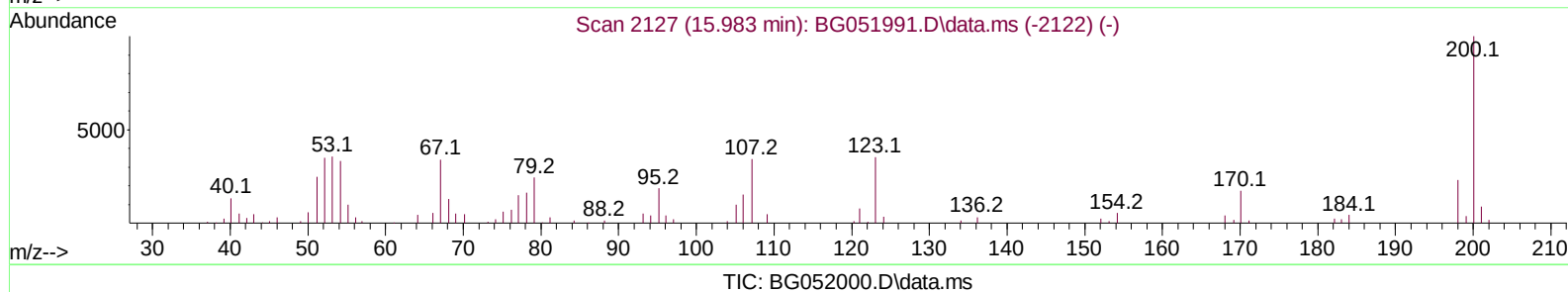
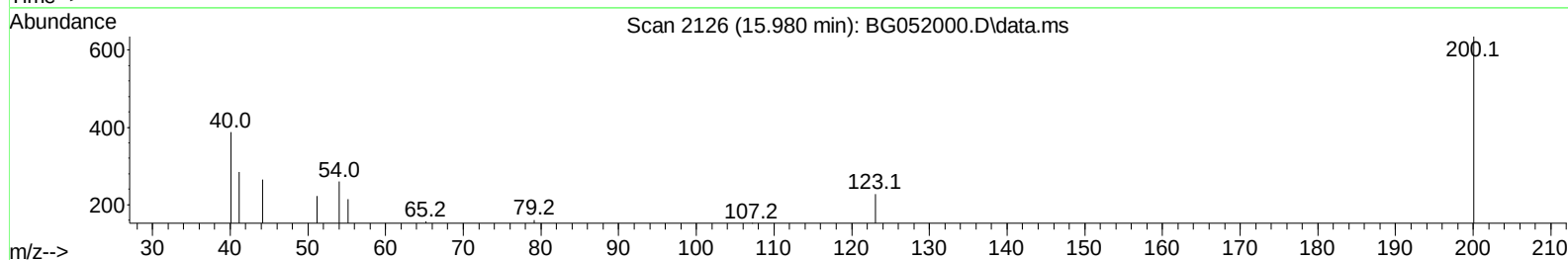
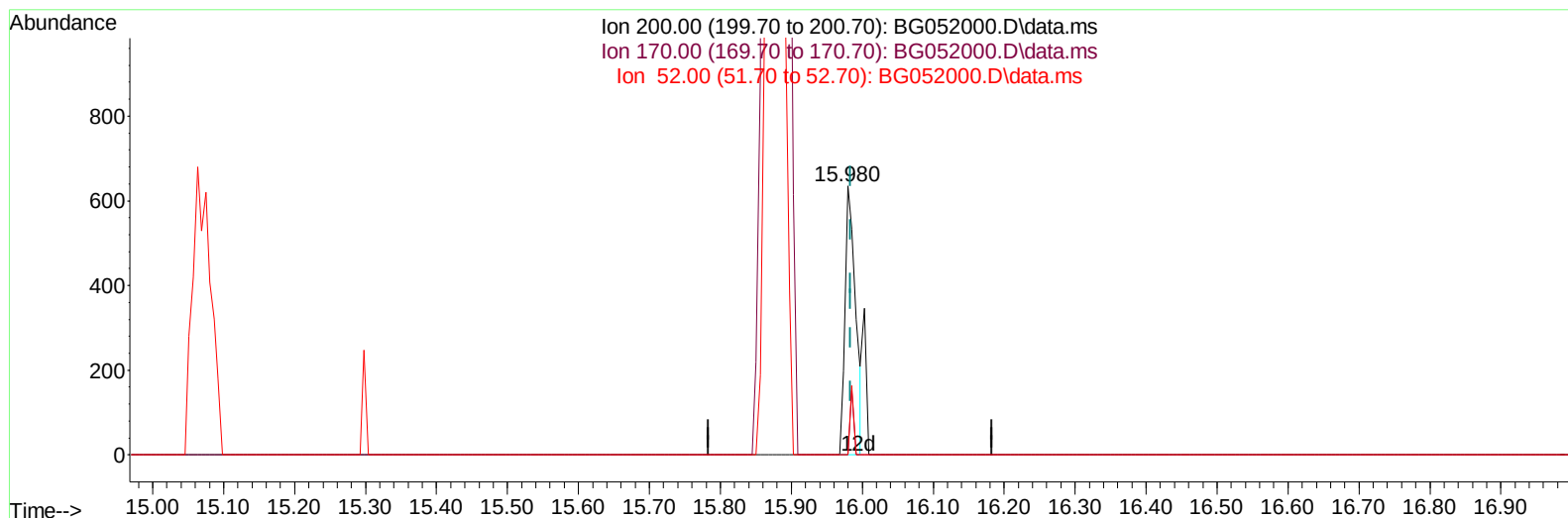
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011522\
 Data File : BG052000.D
 Acq On : 15 Jan 2022 17:05
 Operator : CG/JU
 Sample : N1132-12
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLW3

Manual IntegrationsAPPROVED

Quant Time: Jan 17 04:01:36 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 06 14:43:02 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/17/2022
 Supervised By :mohammad ahmed 01/21/2022



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

15.980min (-0.004) 0.57 ng/ul

response 668

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

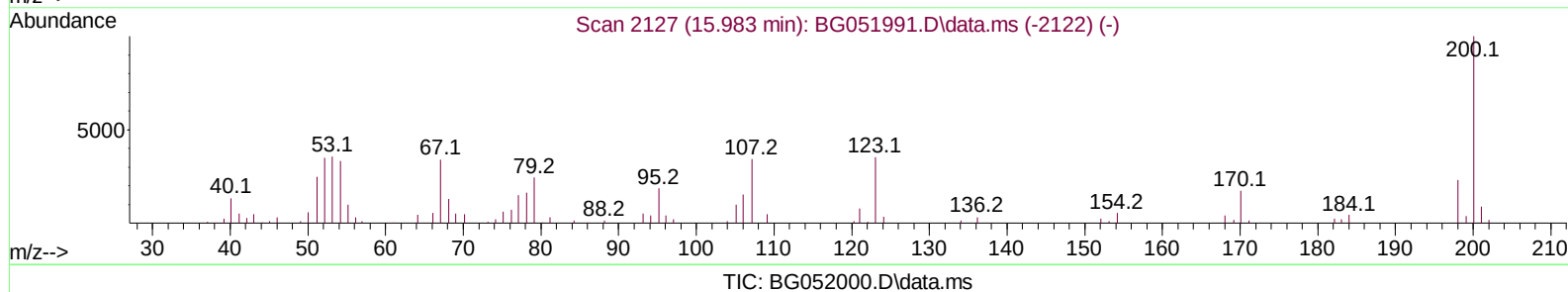
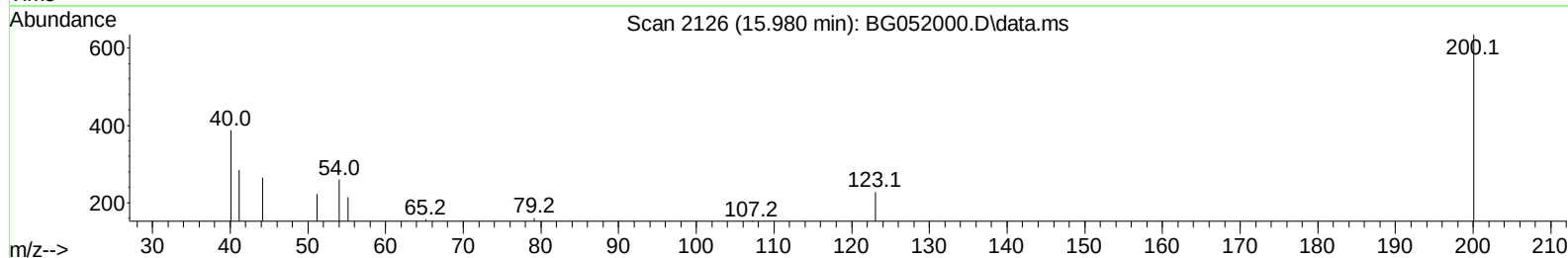
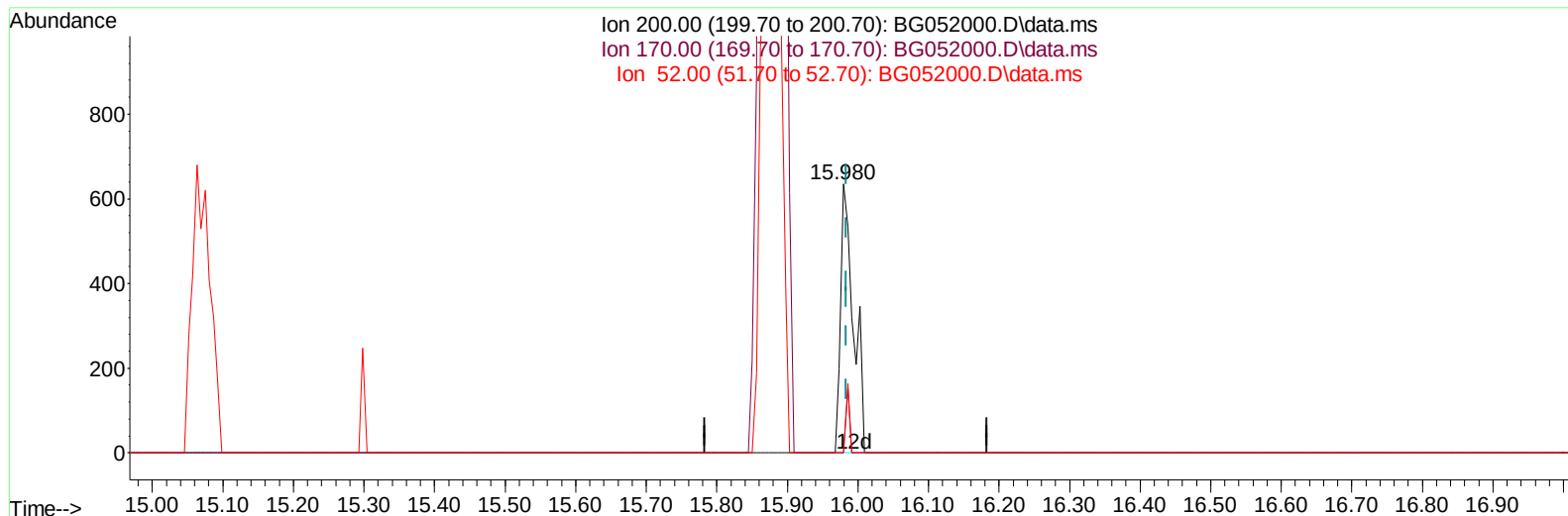
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.980min (-0.004) 0.68 ng/ul m

response 790

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	0.00#
52.00	47.40	0.00#
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	8.249	152	23178	20.000	ng/ul	0.00
20) Naphthalene-d8	11.086	136	101123	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.887	164	74307	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.636	188	184766	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.954	240	162079	20.000	ng/ul	# 0.00
88) Perylene-d12	25.443	264	164160	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.567	96	2772	4.274	ng/uL	-0.01
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.385	99	55015	24.252	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.550	67	39063	27.126	ng/ul	0.00
11) 2-Chlorophenol-d4	7.773	132	36912	24.576	ng/ul	0.00
15) 4-Methylphenol-d8	8.948	113	42496	24.098	ng/ul	0.00
21) Nitrobenzene-d5	9.412	128	22907	28.668	ng/ul	0.00
24) 2-Nitrophenol-d4	10.146	143	14188	16.055	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.698	165	38556	22.321	ng/ul	0.00
31) 4-Chloroaniline-d4	11.210	131	44634	18.554	ng/ul	0.00
46) Dimethylphthalate-d6	14.276	166	157472	25.638	ng/ul	0.00
49) Acenaphthylene-d8	14.581	160	212519	28.644	ng/ul	0.00
54) 4-Nitrophenol-d4	15.063	143	6595	6.422	ng/ul	0.00
60) Fluorene-d10	15.880	176	152249	28.431	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.980	200	790m	0.678	ng/ul	0.00
73) Anthracene-d10	17.736	188	258517	29.495	ng/ul	0.00
81) Pyrene-d10	20.015	212	292106	31.315	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.202	264	246006	28.466	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.409	94	4424	1.911	ng/ul	97

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

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