

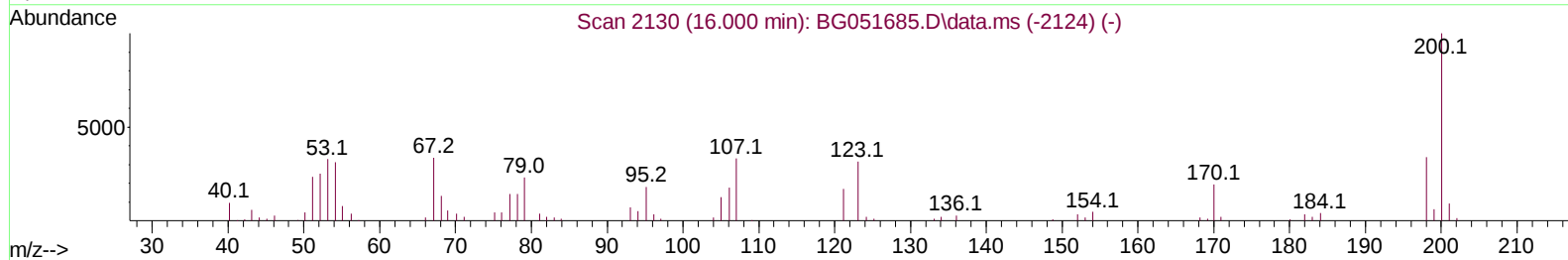
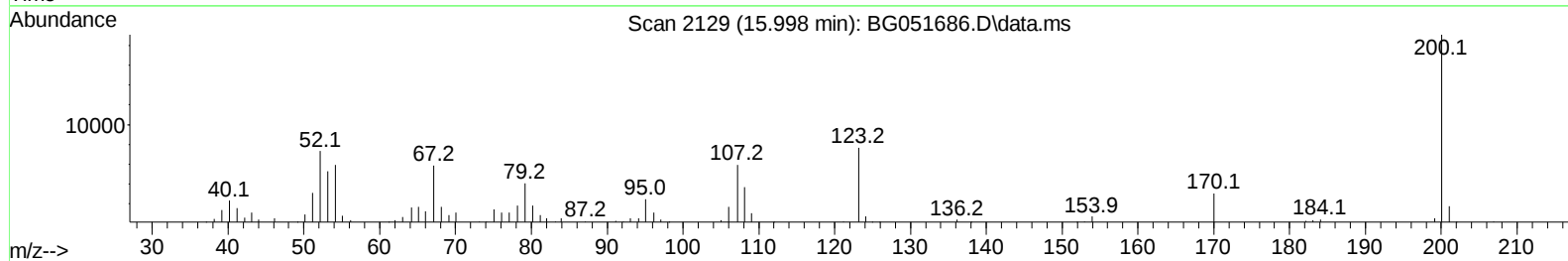
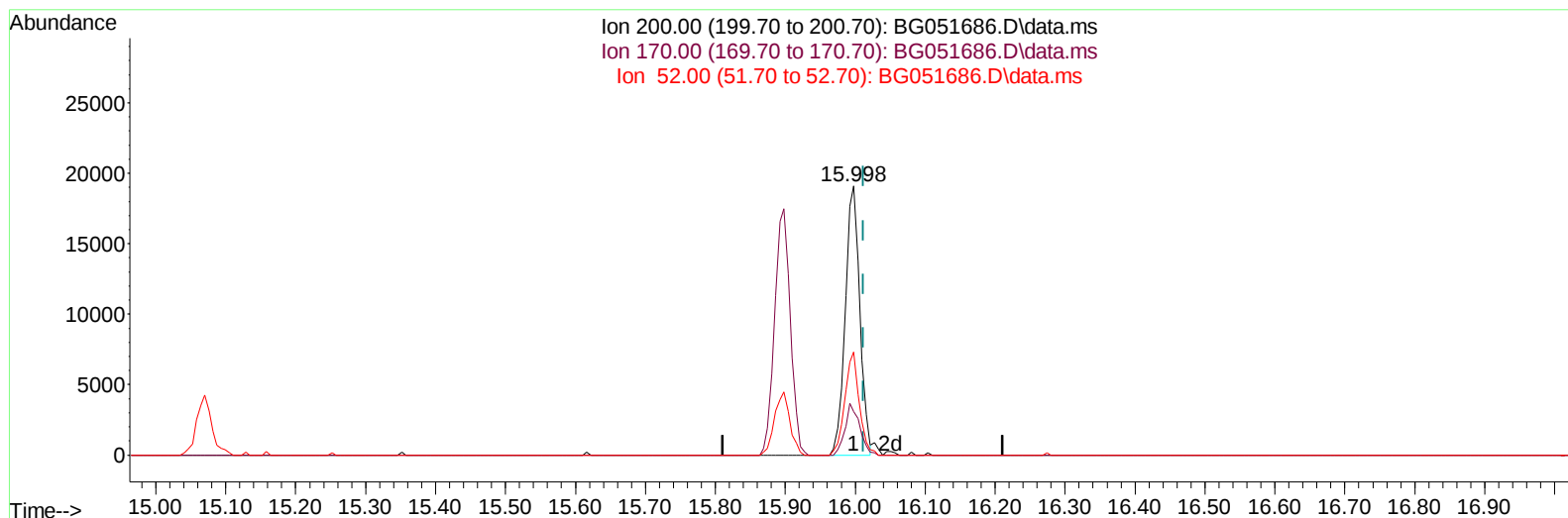
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122021\
Data File : BG051686.D
Acq On : 20 Dec 2021 18:08
Operator : CG/JU
Sample : PB141559BL
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK559

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/21/2021
Supervised By :mohammad ahmed 12/28/2021

Quant Time: Dec 20 23:11:37 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 14 14:19:57 2021
Response via : Initial Calibration



TIC: BG051686.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)**15.998min (-0.013) 21.40 ng/ul****response 27926**

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	16.07
52.00	47.40	38.41
0.00	0.00	0.00

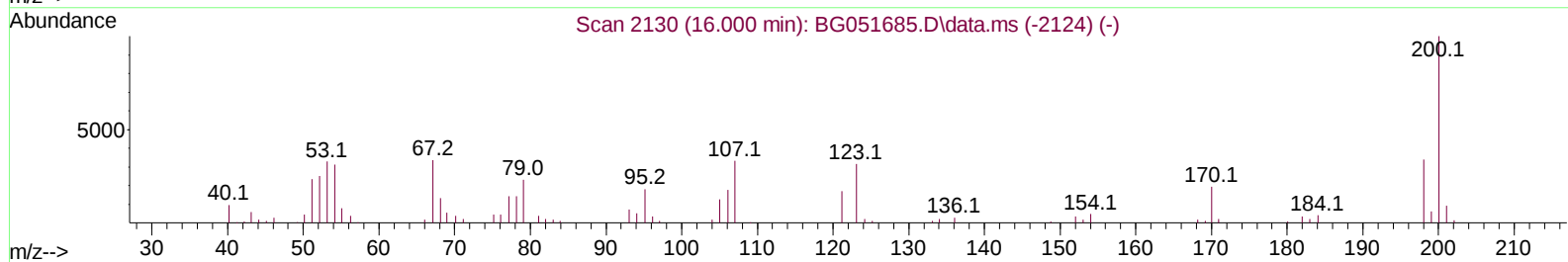
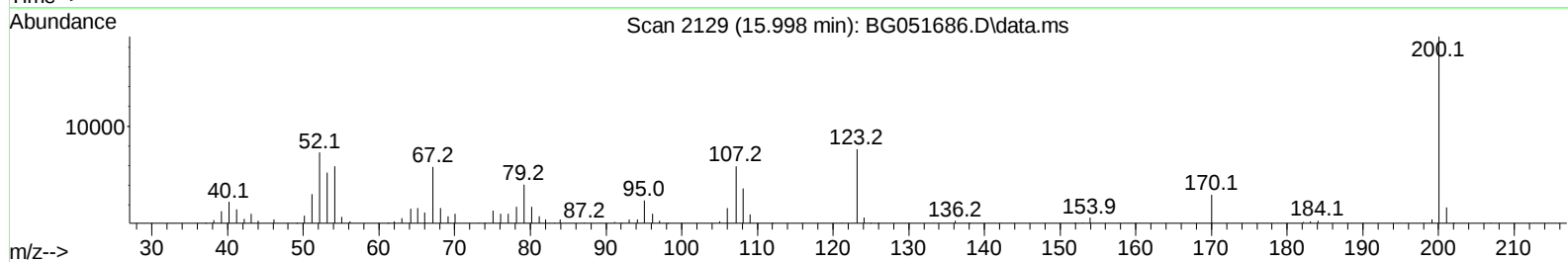
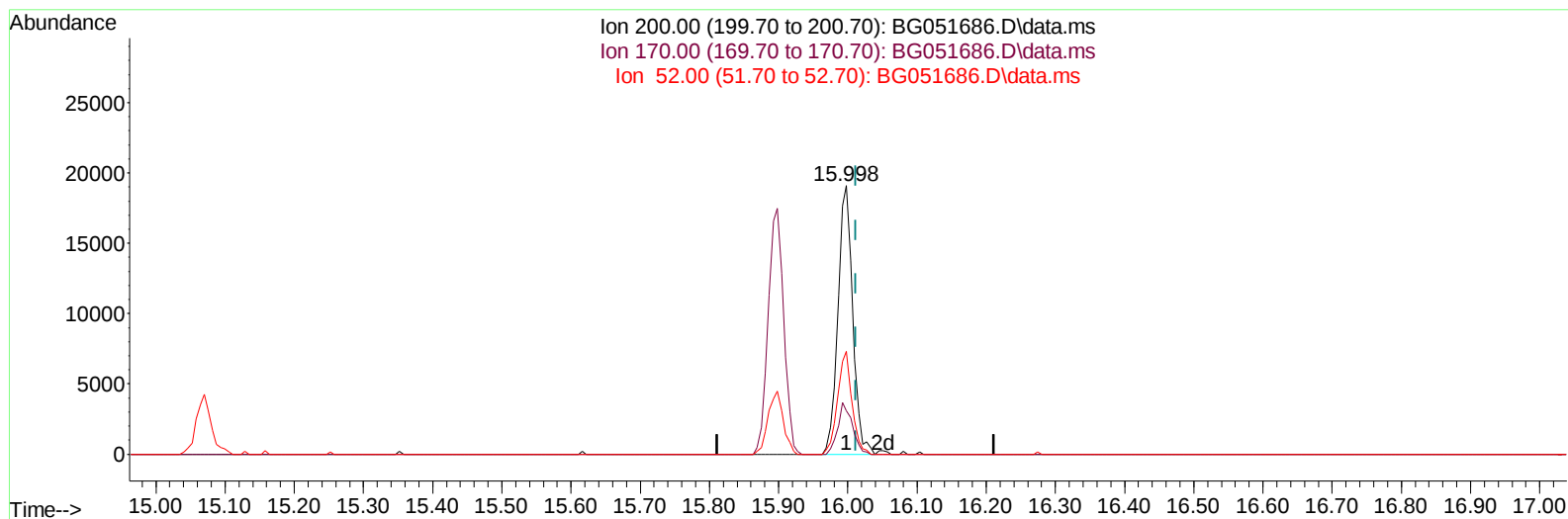
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TIC: BG051686.D\data.ms

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

15.998min (-0.013) 21.75 ng/ul m

response 28393

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	16.07
52.00	47.40	38.41
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	8.273	152	37023	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.111	136	156688	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.911	164	110709	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.655	188	250509	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.972	240	214014	20.000	ng/ul	#-0.01
88) Perylene-d12	25.468	264	214904	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.597	96	4946	5.548	ng/uL	0.00
4) Pyridine-d5	4.026	84	65313	24.248	ng/ul	-0.02
7) Phenol-d5	7.404	99	100881	30.518	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.574	67	58751	29.897	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.797	132	71886	31.267	ng/ul	-0.01
15) 4-Methylphenol-d8	8.966	113	81080	31.731	ng/ul	-0.01
21) Nitrobenzene-d5	9.442	128	37332	32.497	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.171	143	41992	33.438	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.717	165	81426	29.900	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.228	131	91795	25.601	ng/ul	-0.01
46) Dimethylphthalate-d6	14.295	166	266605	30.087	ng/ul	-0.02
49) Acenaphthylene-d8	14.606	160	333610	30.314	ng/ul	0.00
54) 4-Nitrophenol-d4	15.070	143	39504	27.415	ng/ul	0.00
60) Fluorene-d10	15.898	176	229824	28.949	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.998	200	28393m	21.754	ng/ul	-0.01
73) Anthracene-d10	17.755	188	379662	31.771	ng/ul	0.00
81) Pyrene-d10	20.028	212	414917	32.213	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.227	264	363580	32.154	ng/ul	-0.03

Target Compounds Qvalue

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

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