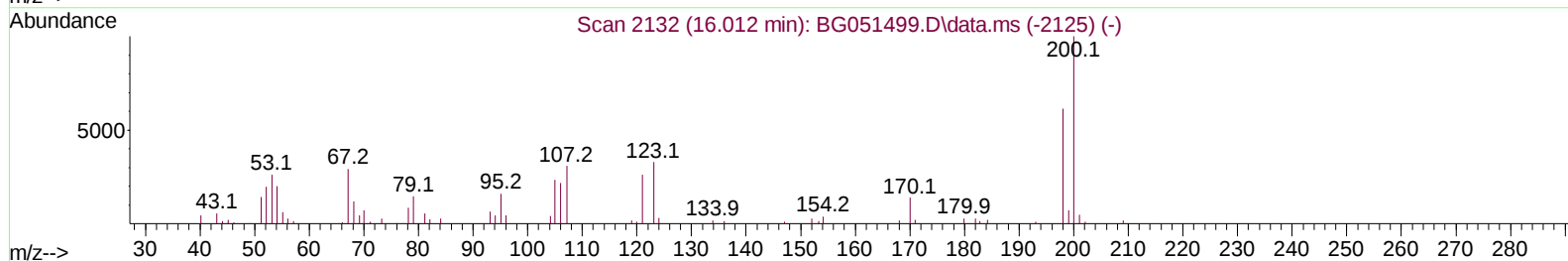
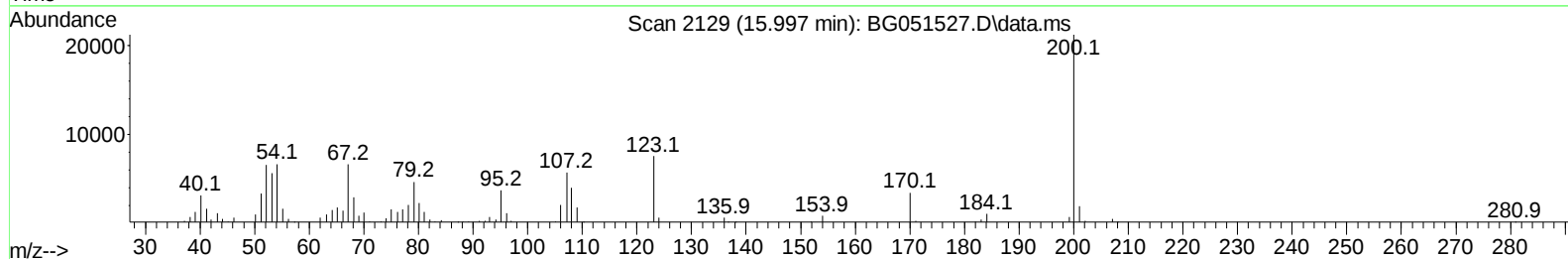
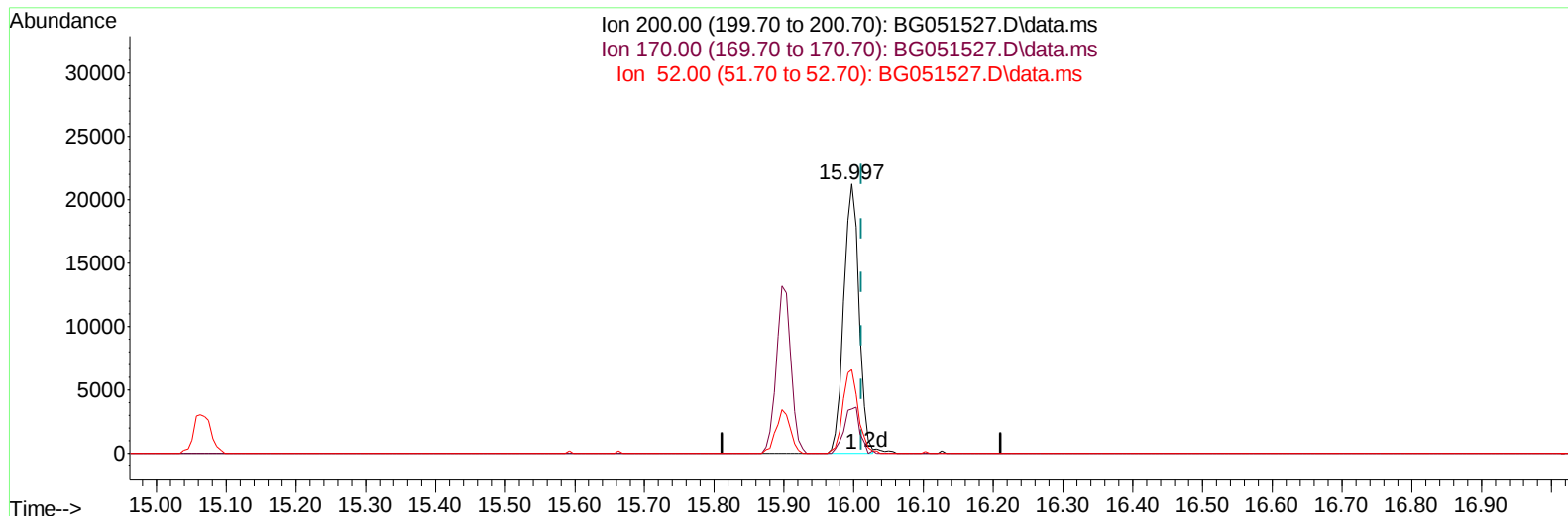


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051527.D
Acq On : 15 Dec 2021 7:27
Operator : CG/JU
Sample : PB141346BL
Misc :
ALS Vial : 25 Sample Multiplier: 1

Quant Time: Dec 15 08:00:51 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: BG051527.D\data.ms

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

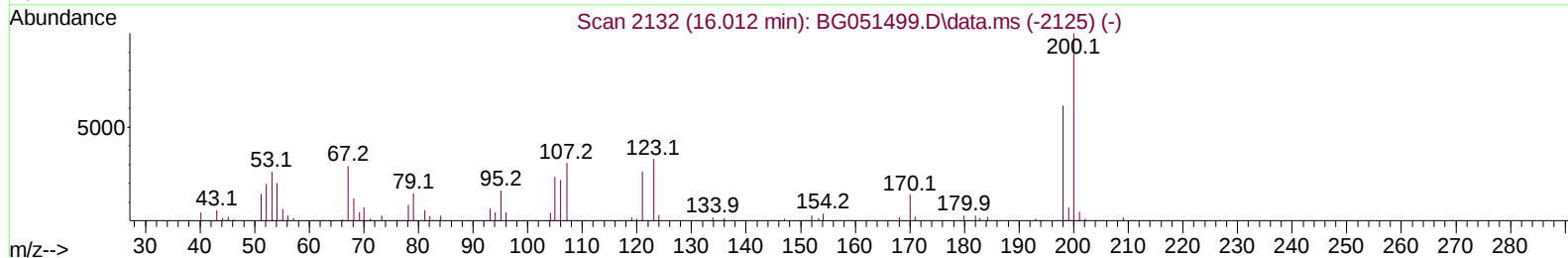
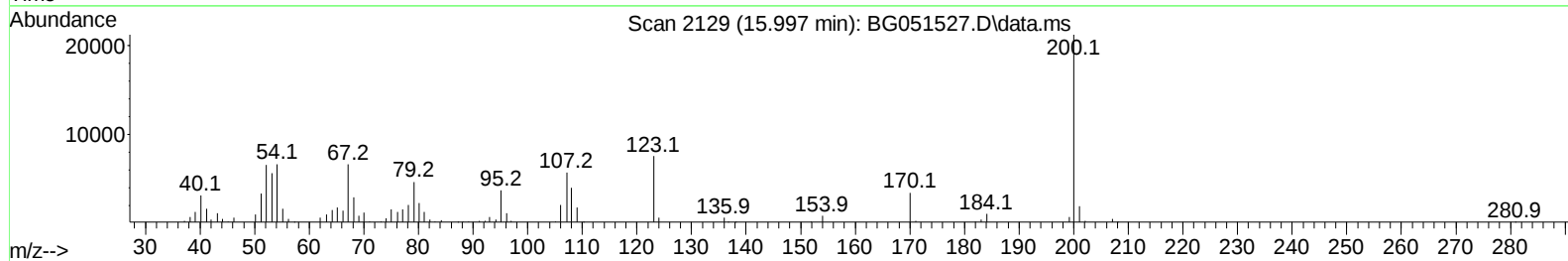
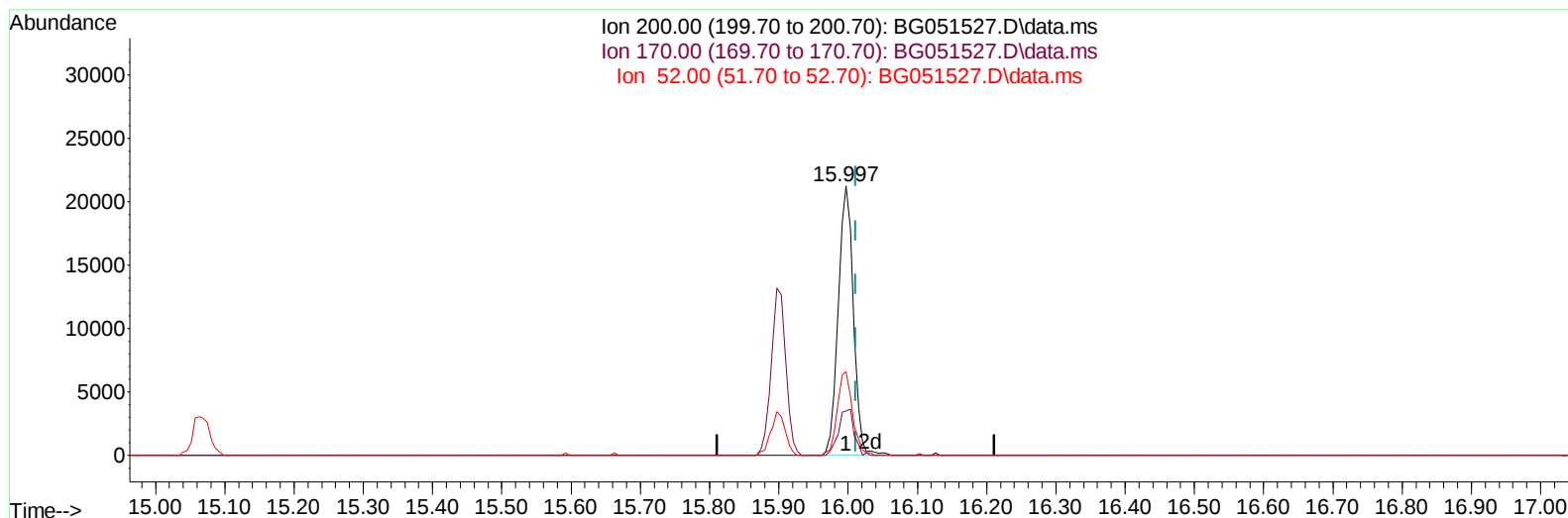
15.997min (-0.014) 31.29 ng/ul

response 31863

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	16.30
52.00	47.40	31.00#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051527.D
Acq On : 15 Dec 2021 7:27
Operator : CG/JU
Sample : PB141346BL
Misc :
ALS Vial : 25 Sample Multiplier: 1

Quant Time: Dec 15 08:00:51 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: BG051527.D\data.ms

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

15.997min (-0.014) 31.53 ng/ul m

response 32101

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	16.30
52.00	47.40	31.00#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
 Data File : BG051527.D
 Acq On : 15 Dec 2021 7:27
 Operator : CG/JU
 Sample : PB141346BL
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Quant Time: Dec 15 08:00:51 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
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.278	152	25723	20.000	ng/ul	0.00
20) Naphthalene-d8	11.116	136	101464	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.910	164	74392	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.654	188	195419	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.977	240	200933	20.000	ng/ul	# 0.00
88) Perylene-d12	25.473	264	200201	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.608	96	3918	6.326	ng/uL	0.00
4) Pyridine-d5	4.031	84	53187	28.421	ng/ul	-0.01
7) Phenol-d5	7.409	99	64359	28.022	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.585	67	40301	29.517	ng/ul	0.00
11) 2-Chlorophenol-d4	7.802	132	46366	29.027	ng/ul	0.00
15) 4-Methylphenol-d8	8.966	113	50471	28.429	ng/ul	-0.01
21) Nitrobenzene-d5	9.447	128	23461	31.538	ng/ul	0.00
24) 2-Nitrophenol-d4	10.176	143	26819	32.980	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.722	165	51590	29.255	ng/ul	0.00
31) 4-Chloroaniline-d4	11.233	131	66262	28.539	ng/ul	0.00
46) Dimethylphthalate-d6	14.305	166	189940	31.900	ng/ul	0.00
49) Acenaphthylene-d8	14.605	160	217672	29.435	ng/ul	0.00
54) 4-Nitrophenol-d4	15.063	143	29543	30.511	ng/ul	-0.01
60) Fluorene-d10	15.897	176	159390	29.878	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.997	200	32101m	31.529	ng/ul	-0.01
73) Anthracene-d10	17.754	188	293973	31.535	ng/ul	0.00
81) Pyrene-d10	20.033	212	375259	31.031	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.232	264	337942	32.082	ng/ul	-0.02

Target Compounds	Qvalue

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

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