

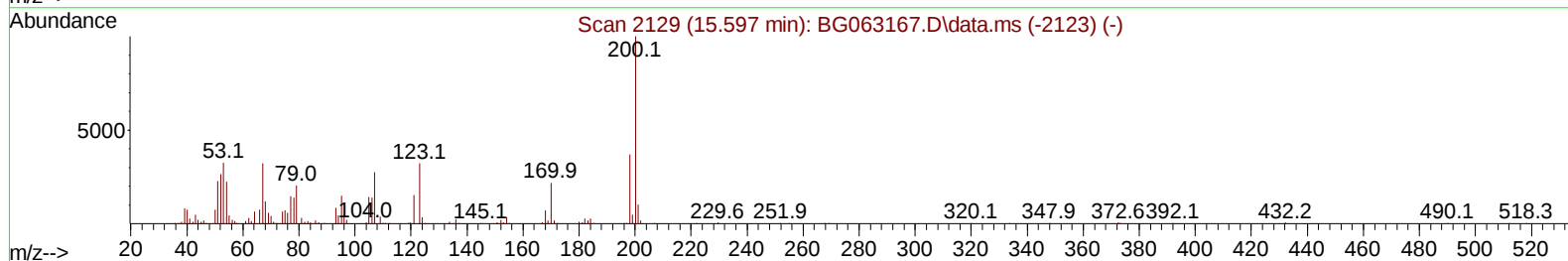
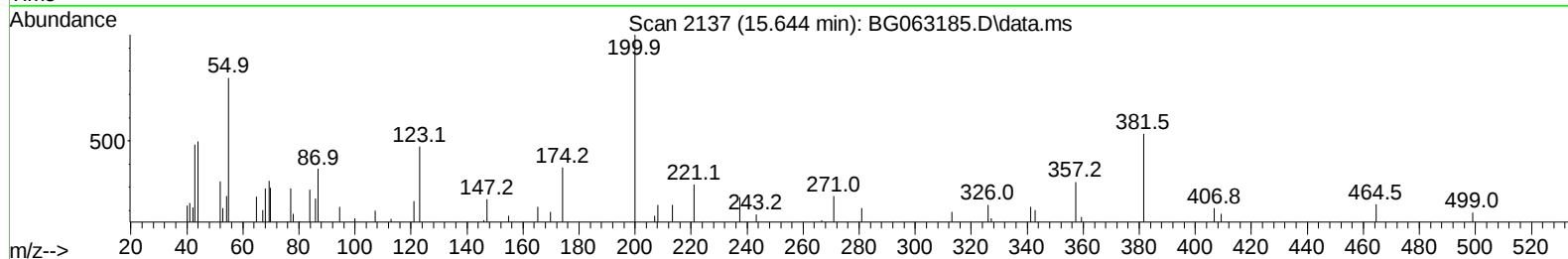
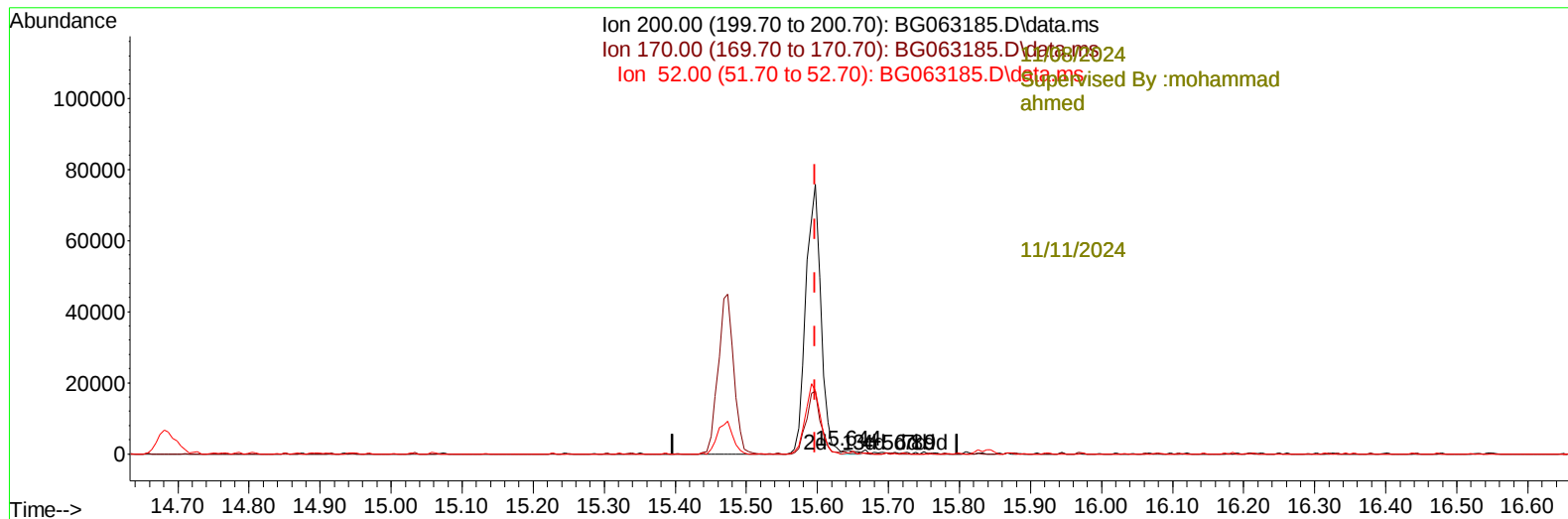
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110624\
Data File : BG063185.D
Acq On : 7 Nov 2024 2:29
Operator : RC/JU
Sample : P4634-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CC0P5

Manual IntegrationsAPPROVED

Quant Time: Nov 07 02:27:36 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 06 15:45:52 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/08/2024
Supervised By :mohammad ahmed 11/11/2024



TIC: BG063185.D\data.ms

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

15.644min (+ 0.048) 0.23 ng/u1

response 1320

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.10	20.27
52.00	29.70	33.96
0.00	0.00	0.00

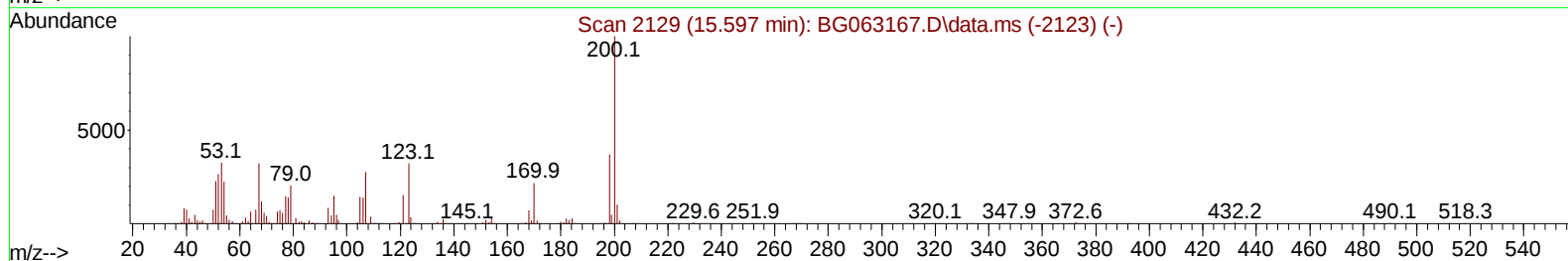
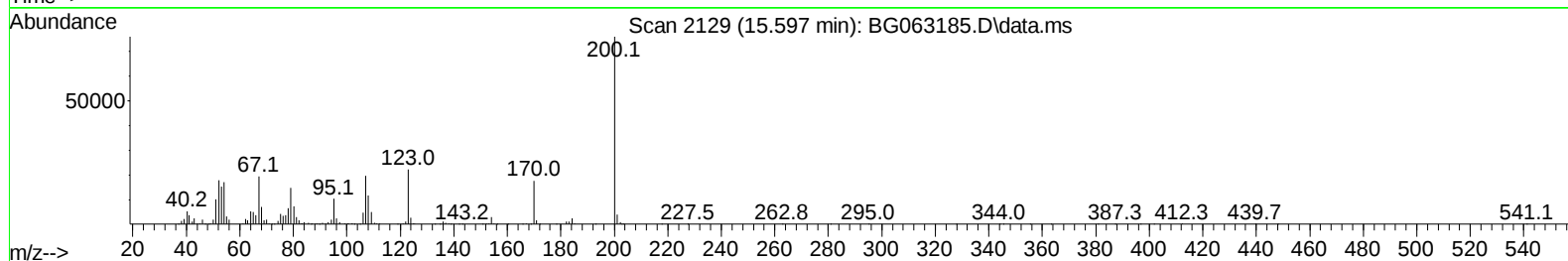
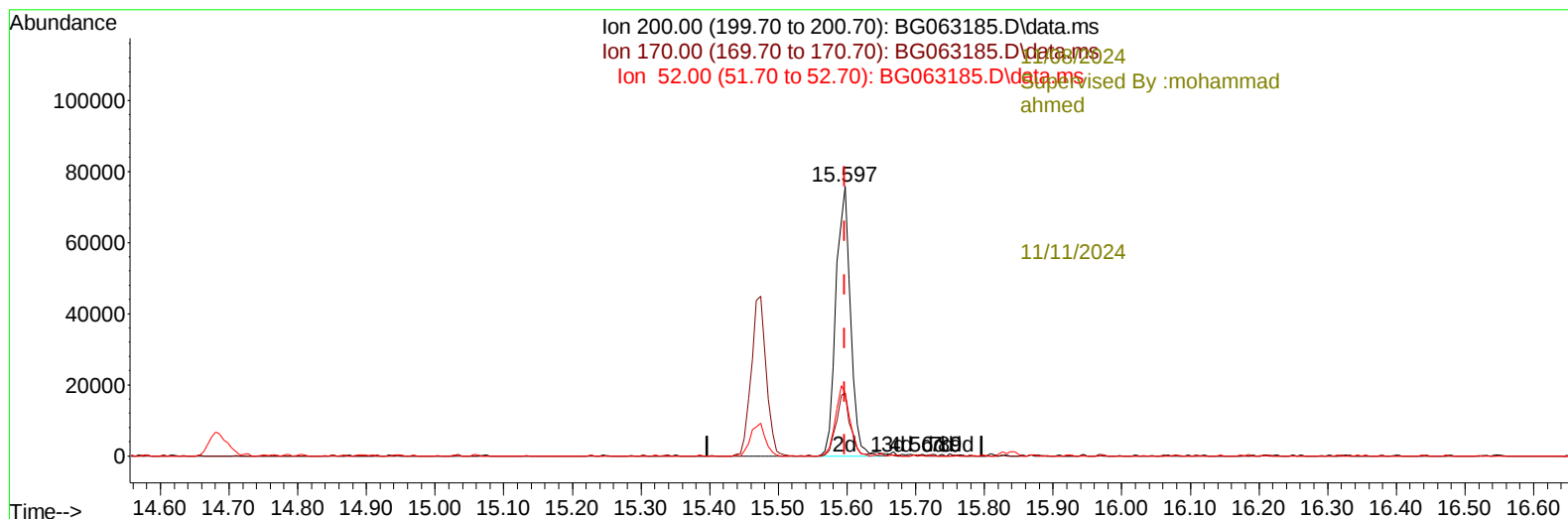
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.597min (+ 0.001) 20.01 ng/ul m

response 112430

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.10	23.30
52.00	29.70	23.46#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/08/2024
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Quant Time: Nov 07 02:28:21 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----11/08/2024						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.842	152	82170	20.000	ng/ul	0.00
20) Naphthalene-d8	10.627	136	381655	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.475	164	357084	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.225	188	894062	20.000	ng/ul	0.00
79) Chrysene-d12	21.479	240	942005	20.000	ng/ul	0.00
88) Perylene-d12	24.510	264	1009540	20.000	ng/ul	0.0011/2024
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.306	96	6533	3.595	ng/uL	0.00
4) Pyridine-d5	3.711	84	88656	14.761	ng/ul	0.00
7) Phenol-d5	7.013	99	166118	22.404	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.172	67	92411	21.248	ng/ul	0.00
11) 2-Chlorophenol-d4	7.372	132	115587	24.018	ng/ul	0.00
15) 4-Methylphenol-d8	8.547	113	141580	22.528	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	60966	22.722	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	79040	22.792	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	164606	24.187	ng/ul	0.00
31) 4-Chloroaniline-d4	10.762	131	189065	21.329	ng/ul	0.00
46) Dimethylphthalate-d6	13.882	166	611643	21.867	ng/ul	0.00
49) Acenaphthylene-d8	14.169	160	668067	24.402	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	74293	19.400	ng/ul	0.00
60) Fluorene-d10	15.468	176	563681	23.878	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.597	200	112430m	20.009	ng/ul	0.00
73) Anthracene-d10	17.325	188	962663	25.053	ng/ul	0.00
81) Pyrene-d10	19.622	212	1216443	25.756	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.311	264	1242321	25.488	ng/ul	0.00

Target Compounds Qvalue

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

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