

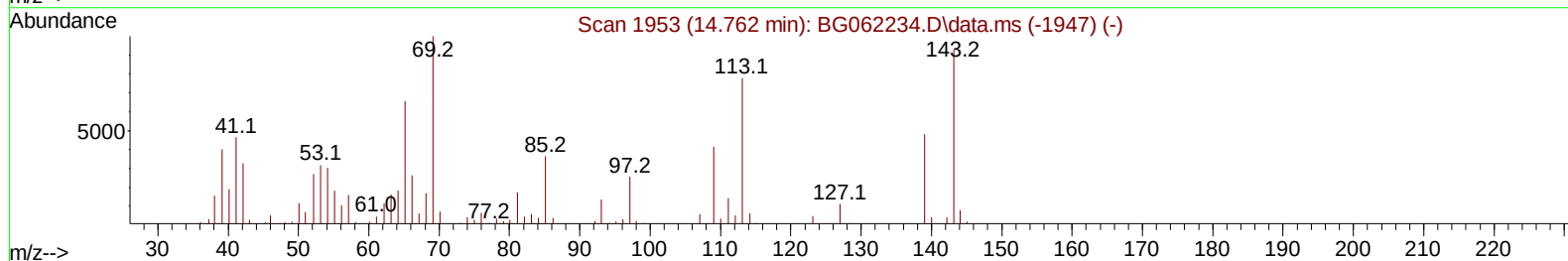
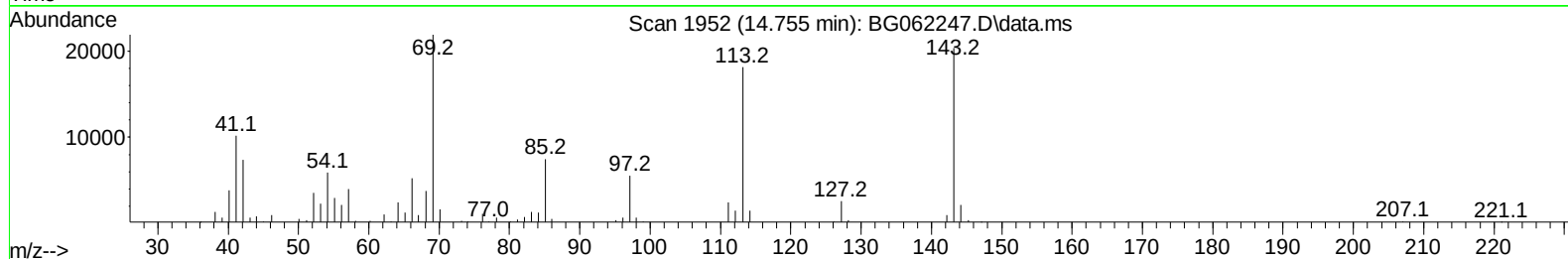
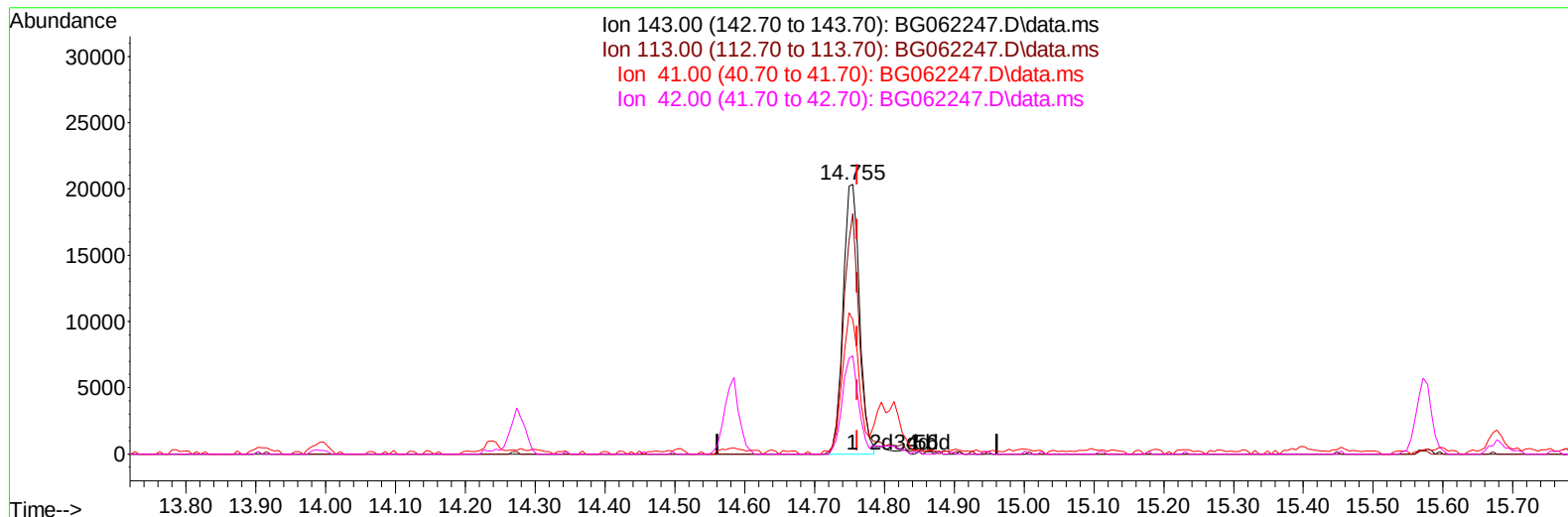
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080524\
Data File : BG062247.D
Acq On : 5 Aug 2024 21:27
Operator : MA/JU
Sample : P3361-08
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E20P0

Manual IntegrationsAPPROVED

Quant Time: Aug 06 02:43:04 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 06 02:21:27 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/06/2024
Supervised By :mohammad ahmed 08/08/2024



TIC: BG062247.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.755min (-0.007) 12.46 ng/ul

response 32769

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	84.90	89.34
41.00	51.00	49.91
42.00	35.30	36.43

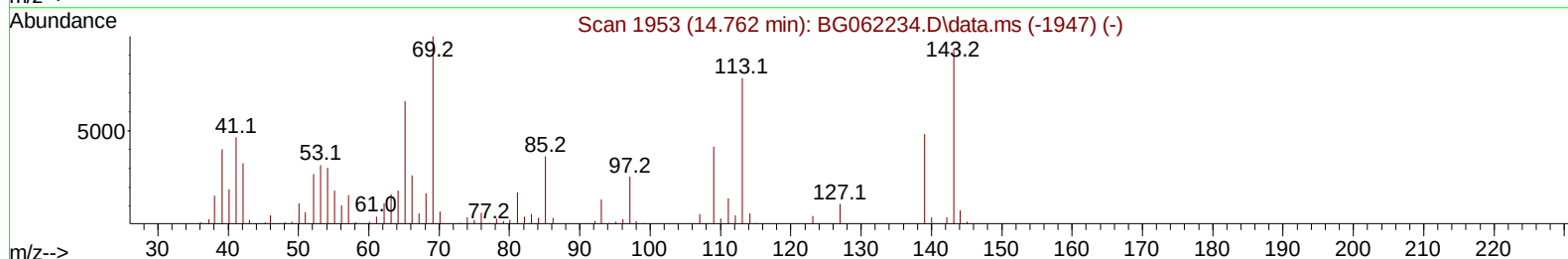
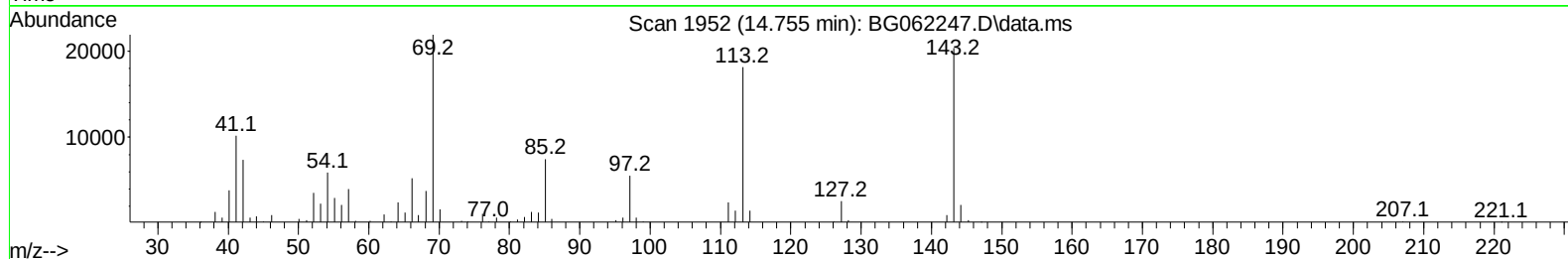
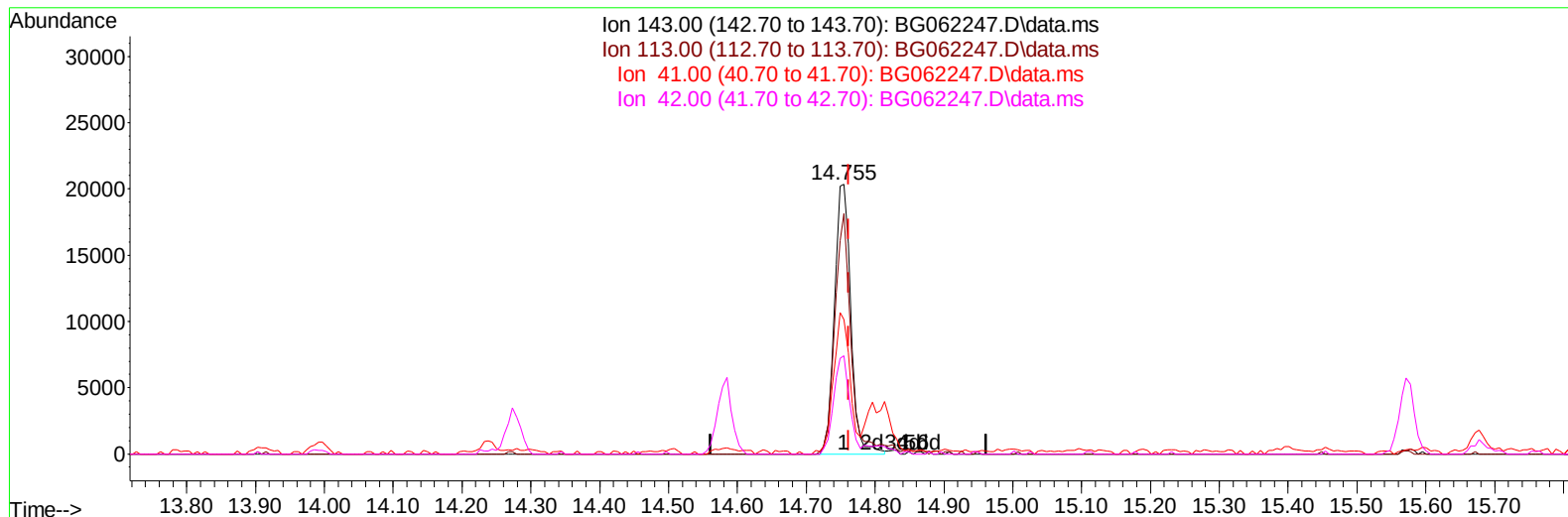
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TIC: BG062247.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.755min (-0.007) 12.73 ng/ul m

response 33487

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	84.90	89.34
41.00	51.00	49.91
42.00	35.30	36.43

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Quant Time: Aug 06 03:07:20 2024
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 QLast Update : Tue Aug 06 02:21:27 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.964	152	67980	20.000	ng/ul	0.00
20) Naphthalene-d8	10.760	136	312336	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.585	164	243553	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	535279	20.000	ng/ul	0.00
79) Chrysene-d12	21.569	240	517397	20.000	ng/ul	0.00
88) Perylene-d12	24.665	264	605896	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.429	96	4569	2.898	ng/uL	0.00
4) Pyridine-d5	3.829	84	54462	10.981	ng/ul	0.00
7) Phenol-d5	7.112	99	106249	16.375	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.294	67	65357	16.363	ng/ul	0.00
11) 2-Chlorophenol-d4	7.494	132	75723	16.289	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	89162	16.498	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	36922	19.171	ng/ul	0.00
24) 2-Nitrophenol-d4	9.838	143	35581	19.639	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.373	165	90565	17.211	ng/ul	0.00
31) 4-Chloroaniline-d4	10.884	131	110588	14.490	ng/ul	0.00
46) Dimethylphthalate-d6	13.991	166	303274	16.139	ng/ul	0.00
49) Acenaphthylene-d8	14.273	160	346221	16.250	ng/ul	0.00
54) 4-Nitrophenol-d4	14.755	143	33487m	12.729	ng/ul	0.00
60) Fluorene-d10	15.572	176	303369	17.986	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.677	200	21433	13.520	ng/ul	0.00
73) Anthracene-d10	17.422	188	454496	17.604	ng/ul	0.00
81) Pyrene-d10	19.701	212	580192	18.866	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.454	264	593423	18.321	ng/ul	-0.01

Target Compounds Qvalue

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

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