

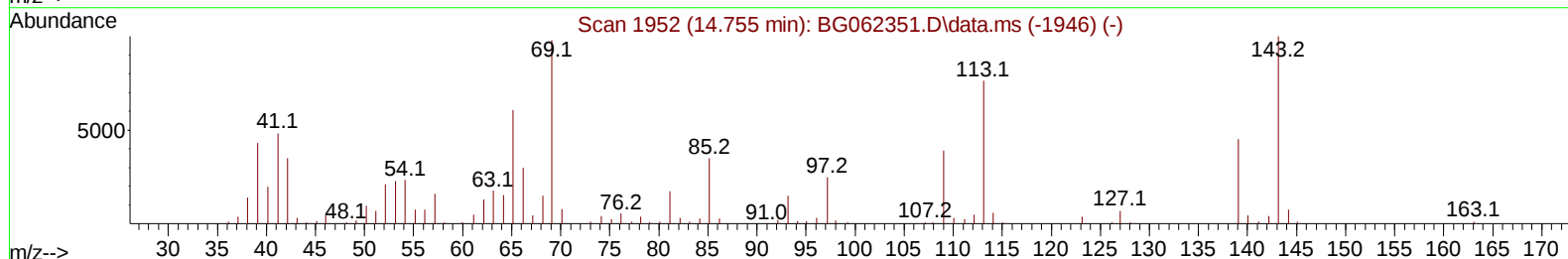
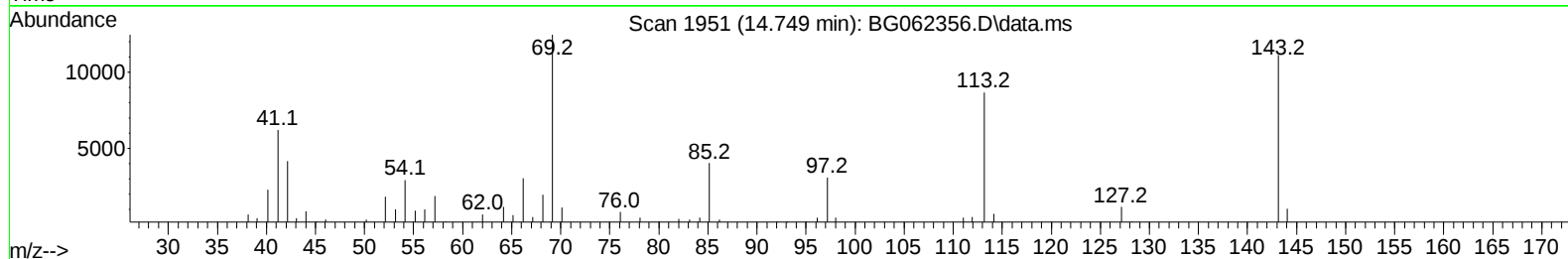
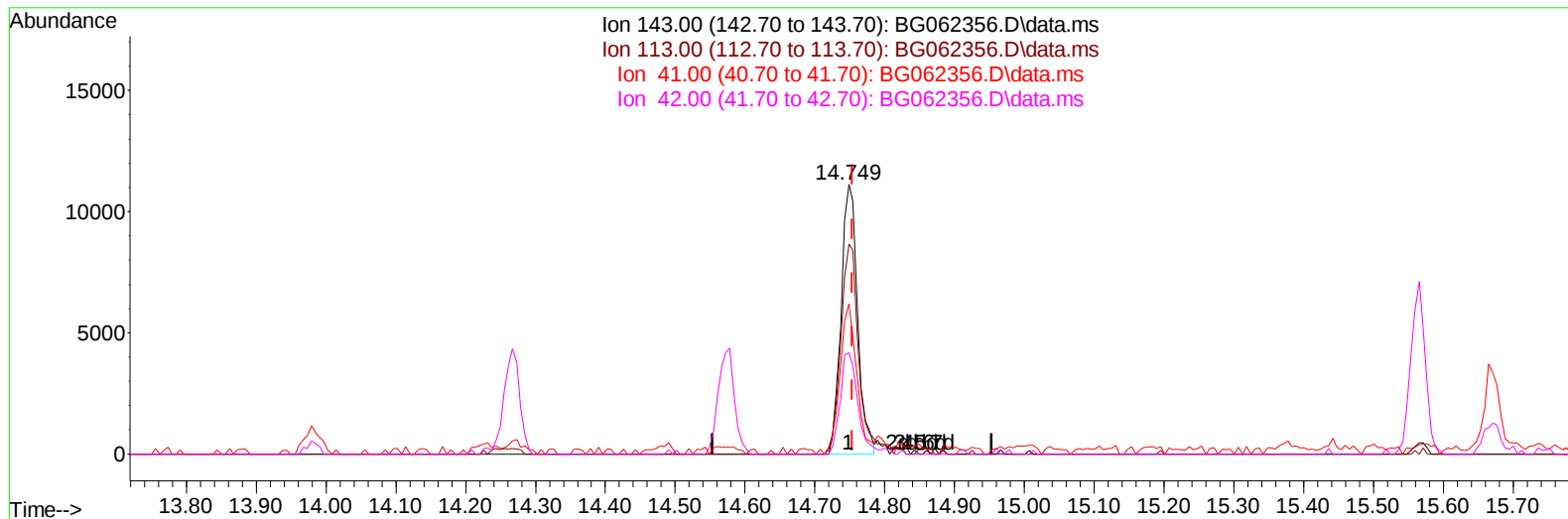
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080824\
Data File : BG062356.D
Acq On : 9 Aug 2024 11:53
Operator : MA/JU
Sample : P3480-18
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E0AB5

Manual IntegrationsAPPROVED

Quant Time: Aug 09 12:37:23 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Aug 09 01:00:03 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/12/2024
Supervised By :mohammad ahmed 08/12/2024



TIC: BG062356.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.749min (-0.005) 8.21 ng/u1

response 18090

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	84.90	78.02
41.00	51.00	55.88
42.00	35.30	37.57

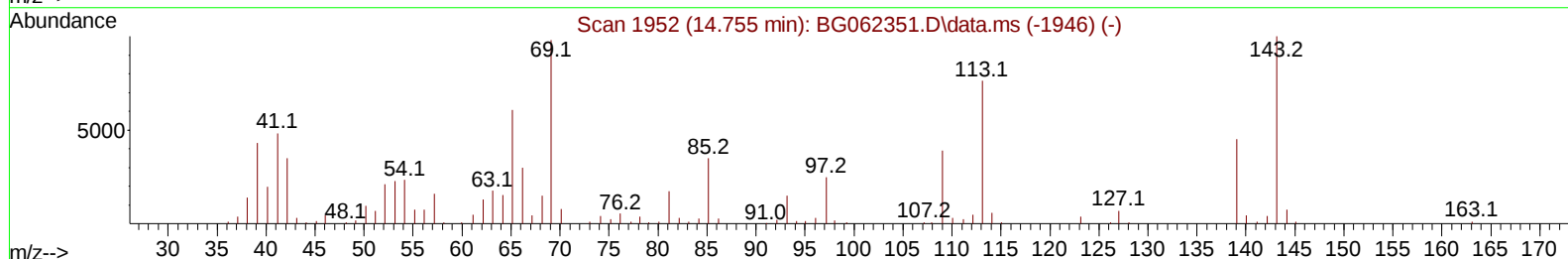
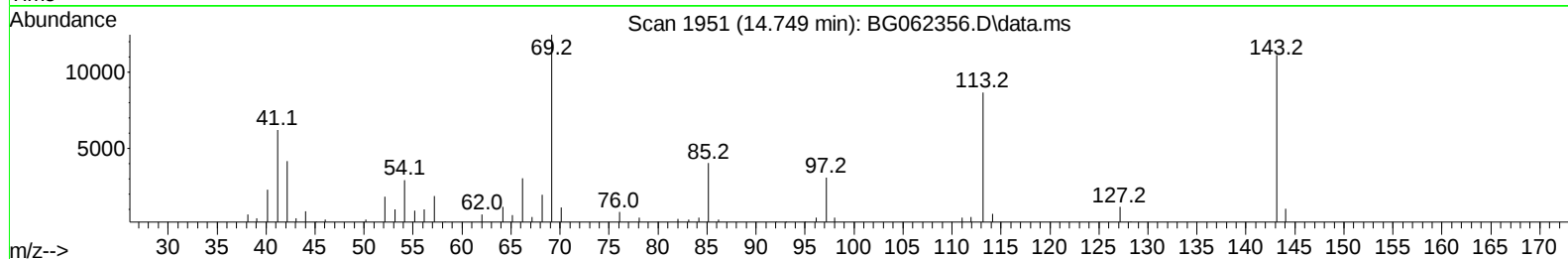
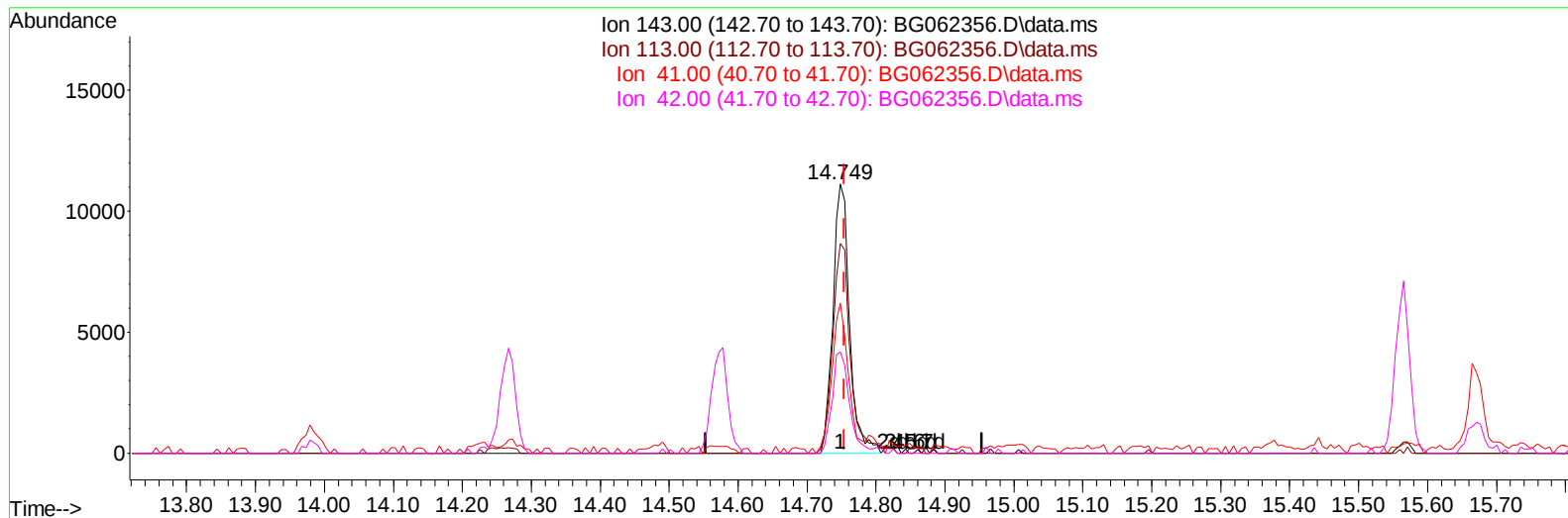
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TIC: BG062356.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.749min (-0.005) 8.41 ng/u1 m

response 18538

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	84.90	78.02
41.00	51.00	55.88
42.00	35.30	37.57

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	53372	20.000	ng/ul	0.00
20) Naphthalene-d8	10.748	136	261088	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.572	164	204044	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.310	188	439110	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	429549	20.000	ng/ul	0.00
88) Perylene-d12	24.635	264	500316	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.423	96	5328	4.304	ng/uL	0.00
4) Pyridine-d5	3.828	84	18799	4.828	ng/ul	0.00
7) Phenol-d5	7.112	99	49101	9.639	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.282	67	82119	26.187	ng/ul	0.00
11) 2-Chlorophenol-d4	7.488	132	95127	26.063	ng/ul	0.00
15) 4-Methylphenol-d8	8.651	113	83682	19.722	ng/ul	0.00
21) Nitrobenzene-d5	9.103	128	51714	32.122	ng/ul	0.00
24) 2-Nitrophenol-d4	9.832	143	58029	38.317	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.366	165	125456	28.521	ng/ul	0.00
31) 4-Chloroaniline-d4	10.877	131	128575	20.153	ng/ul	0.00
46) Dimethylphthalate-d6	13.979	166	412967	26.232	ng/ul	0.00
49) Acenaphthylene-d8	14.267	160	457557	25.634	ng/ul	0.00
54) 4-Nitrophenol-d4	14.749	143	18538m	8.411	ng/ul	0.00
60) Fluorene-d10	15.565	176	365972	25.899	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.671	200	47662	36.650	ng/ul	0.00
73) Anthracene-d10	17.410	188	561533	26.513	ng/ul	0.00
81) Pyrene-d10	19.695	212	701597	27.480	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.430	264	736848	27.550	ng/ul	0.00

Target Compounds Qvalue

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

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