

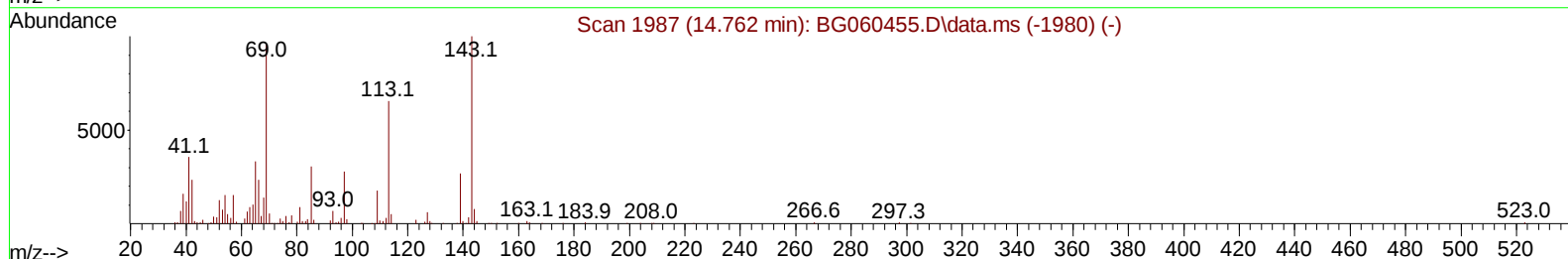
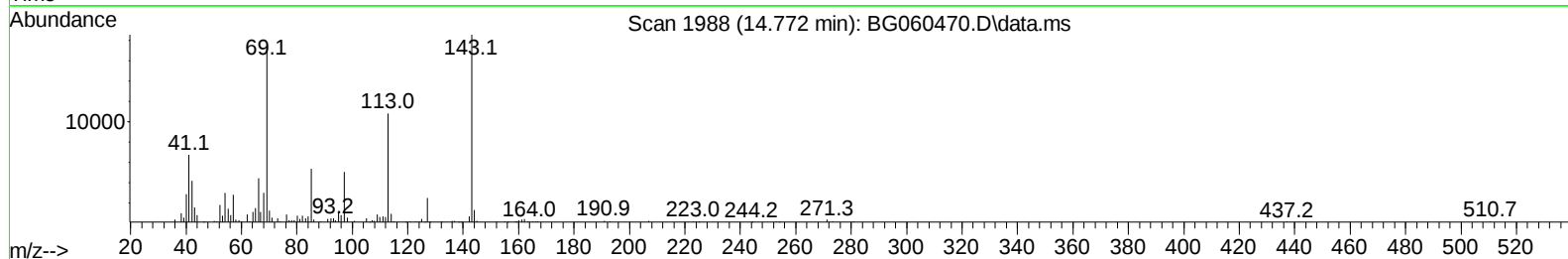
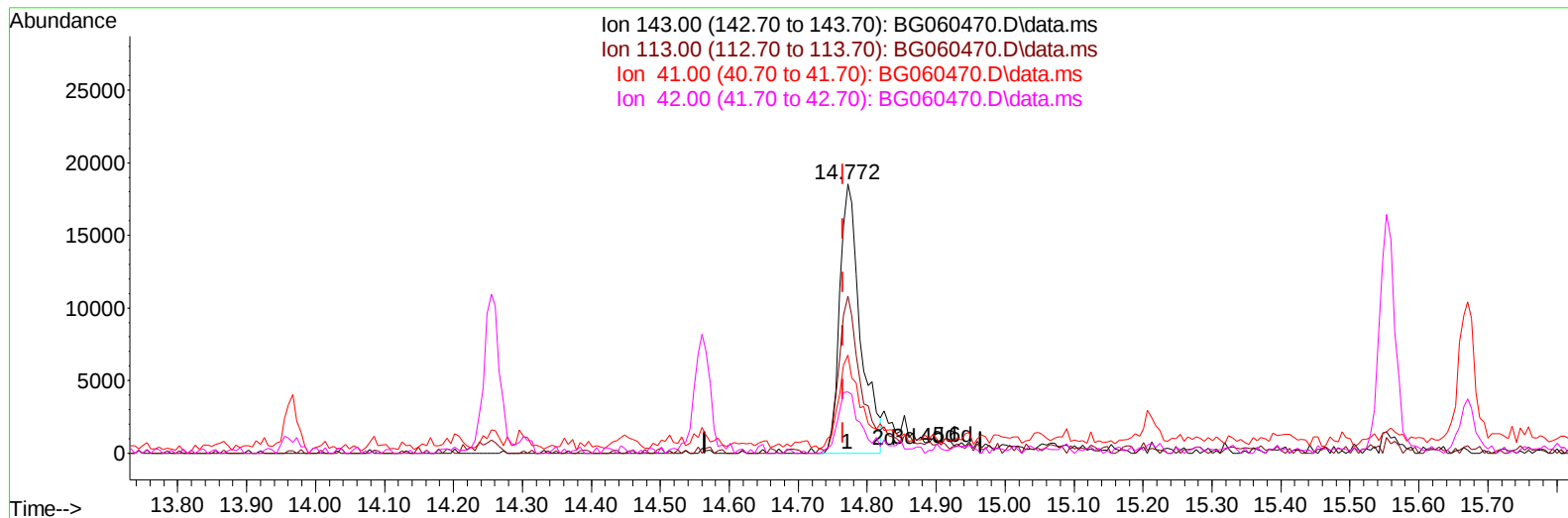
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012924\
Data File : BG060470.D
Acq On : 29 Jan 2024 20:21
Operator : MA/JU
Sample : P1239-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0FC5

Manual IntegrationsAPPROVED

Quant Time: Jan 30 01:49:06 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jan 26 06:06:34 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/30/2024
Supervised By :mohammad ahmed 01/31/2024



TIC: BG060470.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.772min (+ 0.007) 4.72 ng/u1

response 38564

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	58.39
41.00	29.80	36.46#
42.00	22.70	22.66

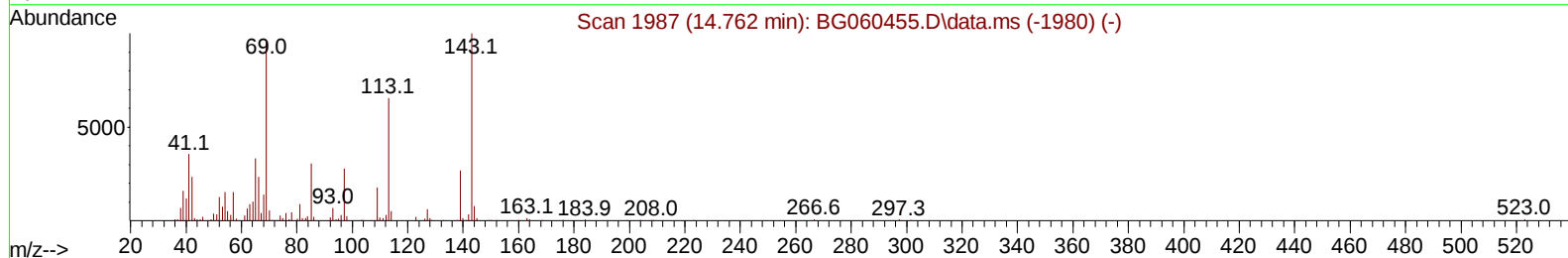
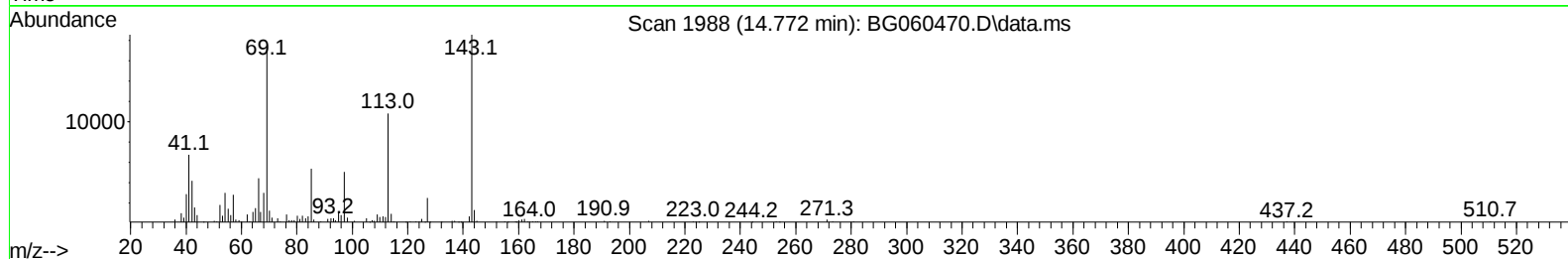
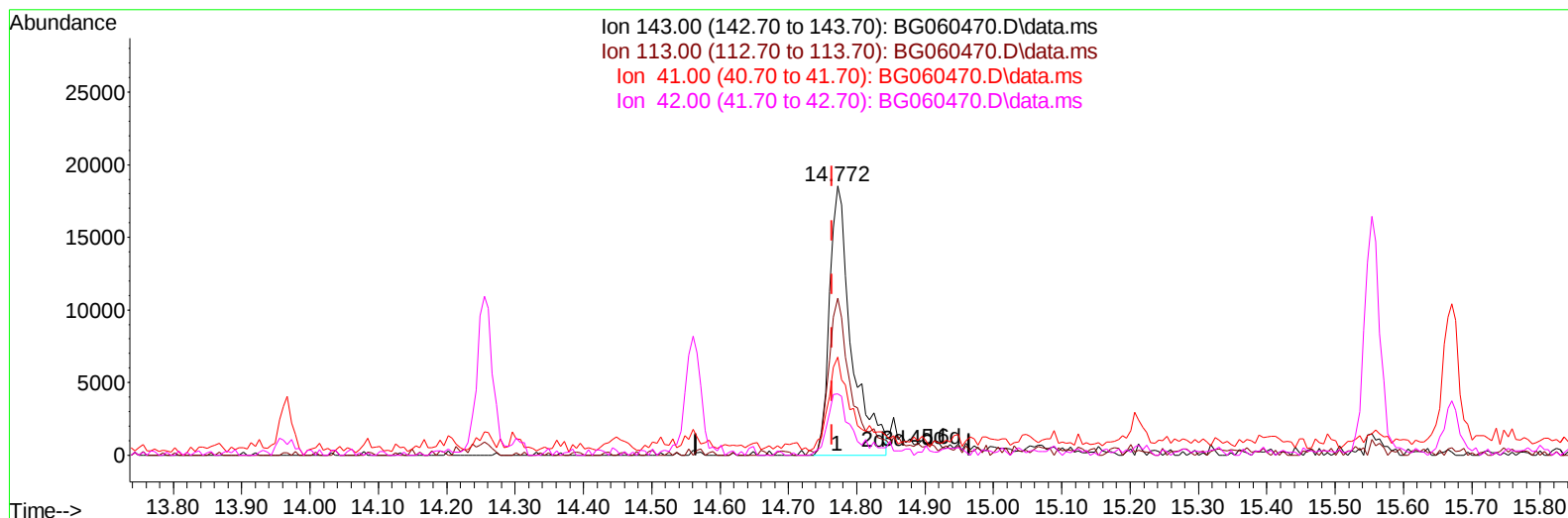
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TIC: BG060470.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.772min (+ 0.007) 5.08 ng/ul m

response 41524

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	58.39
41.00	29.80	36.46#
42.00	22.70	22.66

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/30/2024
 Supervised By :mohammad ahmed 01/31/2024

Quant Time: Jan 30 02:02:49 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 26 06:06:34 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.927	152	206142	20.000	ng/ul	0.00
20) Naphthalene-d8	10.729	136	933302	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.560	164	697408	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.310	188	1769729	20.000	ng/ul	0.00
79) Chrysene-d12	21.576	240	1617416	20.000	ng/ul	0.00
88) Perylene-d12	24.737	264	1812766	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.362	96	19324	4.697	ng/uL	0.00
4) Pyridine-d5	3.773	84	127419	9.791	ng/ul	0.00
7) Phenol-d5	7.087	99	115912	5.921	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.251	67	344035	27.964	ng/ul	0.00
11) 2-Chlorophenol-d4	7.457	132	298910	22.863	ng/ul	0.00
15) 4-Methylphenol-d8	8.632	113	237455	14.725	ng/ul	0.00
21) Nitrobenzene-d5	9.084	128	212101	30.807	ng/ul	0.00
24) 2-Nitrophenol-d4	9.807	143	244748	30.427	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.348	165	509486	30.457	ng/ul	0.00
31) 4-Chloroaniline-d4	10.865	131	465575	21.216	ng/ul	0.00
46) Dimethylphthalate-d6	13.967	166	1885763	33.044	ng/ul	0.00
49) Acenaphthylene-d8	14.255	160	2130716	33.285	ng/ul	0.00
54) 4-Nitrophenol-d4	14.772	143	41524m	5.080	ng/ul	0.00
60) Fluorene-d10	15.553	176	1685954	33.717	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.671	200	296845	27.088	ng/ul	0.00
73) Anthracene-d10	17.410	188	2727374	32.511	ng/ul	0.00
81) Pyrene-d10	19.701	212	3173406	32.736	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.519	264	3184855	33.492	ng/ul	0.00

Target Compounds Qvalue

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

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