

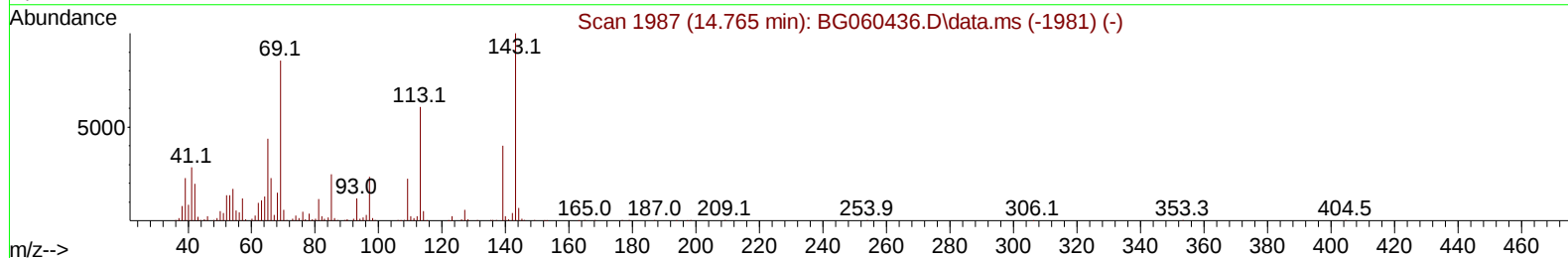
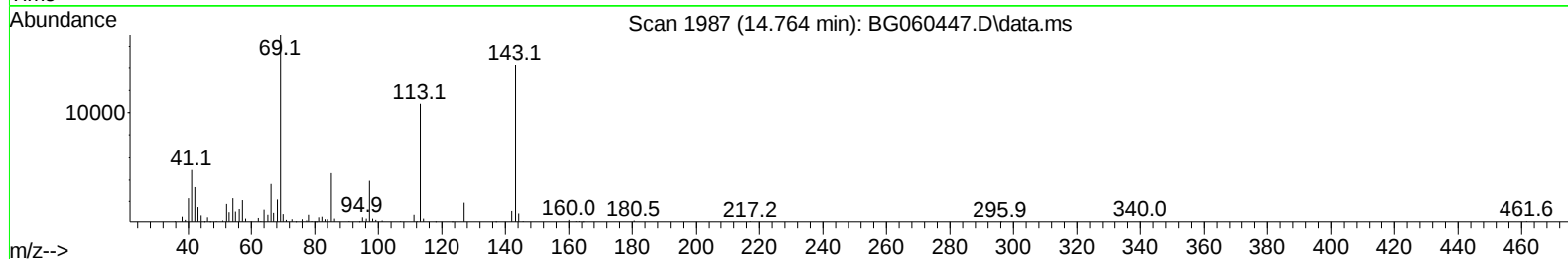
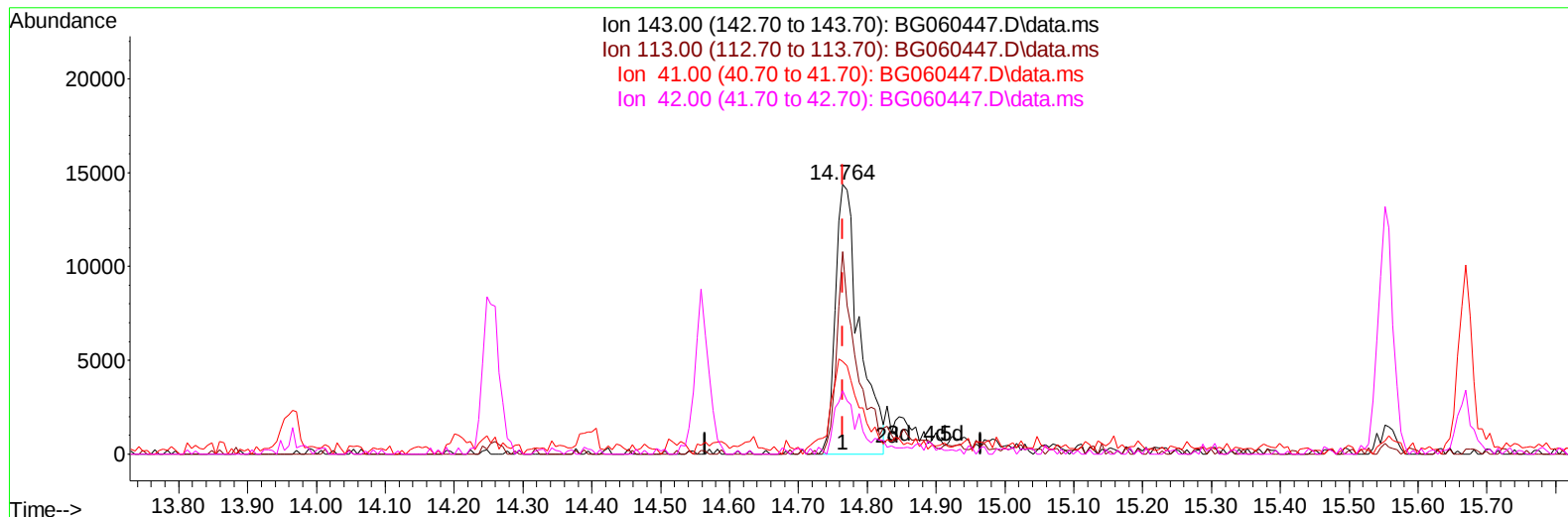
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012624\
Data File : BG060447.D
Acq On : 26 Jan 2024 18:58
Operator : MA/JU
Sample : P1239-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0FB9

Manual IntegrationsAPPROVED

Quant Time: Jan 26 22:04:55 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 01/29/2024
Supervised By :mohammad ahmed 01/30/2024



TIC: BG060447.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.764min (-0.000) 5.01 ng/u1

response 34724

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	75.17#
41.00	29.80	34.25
42.00	22.70	23.49

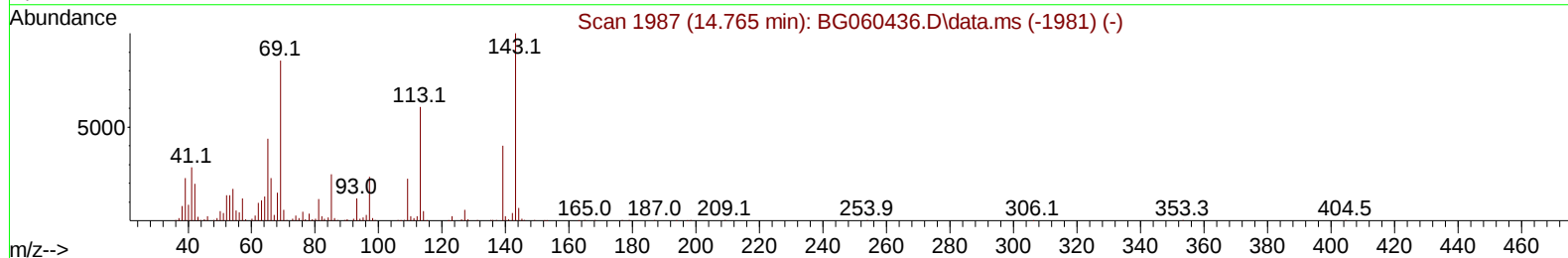
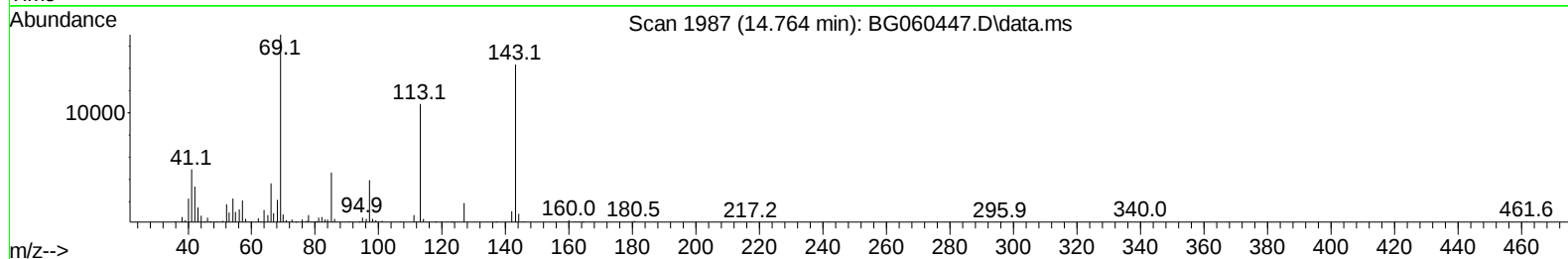
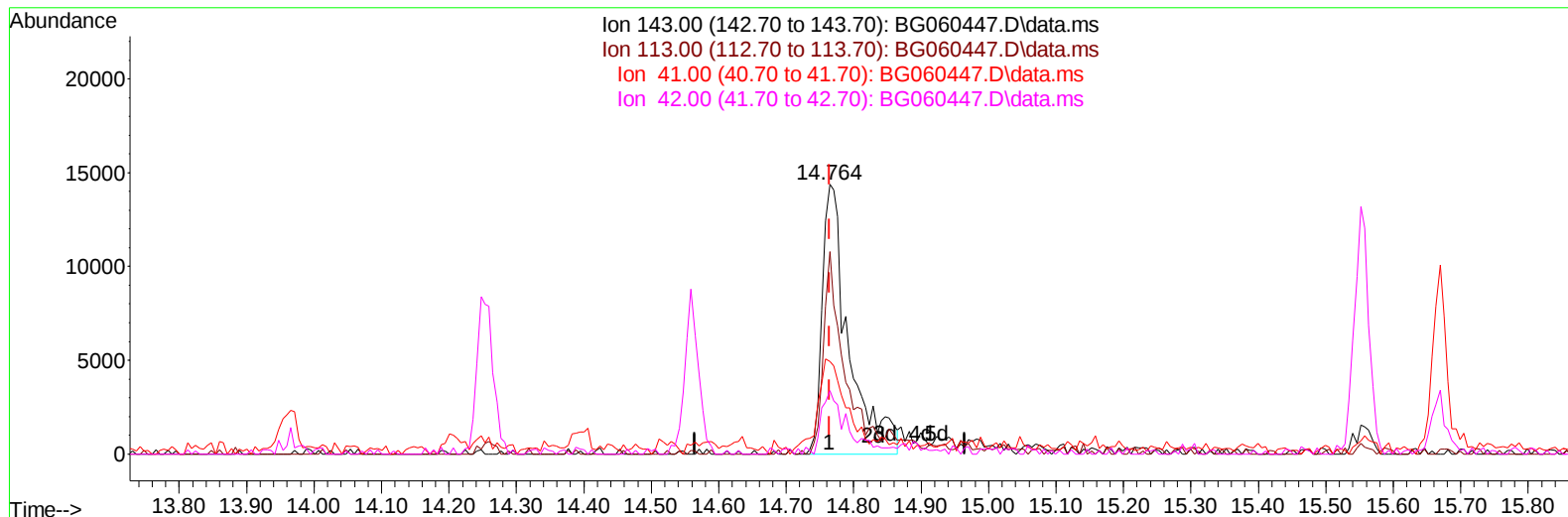
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TIC: BG060447.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.764min (-0.000) 5.64 ng/ul m

response 39042

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	75.17#
41.00	29.80	34.25
42.00	22.70	23.49

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 ALS Vial : 13 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/29/2024
 Supervised By :mohammad ahmed 01/30/2024

Quant Time: Jan 26 22:15:14 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	7.925	152	163734	20.000	ng/ul	0.00
20) Naphthalene-d8	10.728	136	749501	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.559	164	591107	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.309	188	1470505	20.000	ng/ul	0.00
79) Chrysene-d12	21.568	240	1420632	20.000	ng/ul	0.00
88) Perylene-d12	24.729	264	1590150	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.360	96	18958	5.802	ng/uL	0.00
4) Pyridine-d5	3.777	84	84415	8.167	ng/ul	0.00
7) Phenol-d5	7.091	99	93309	6.000	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.250	67	263858	27.002	ng/ul	0.00
11) 2-Chlorophenol-d4	7.455	132	220492	21.233	ng/ul	0.00
15) 4-Methylphenol-d8	8.630	113	179510	14.014	ng/ul	0.00
21) Nitrobenzene-d5	9.083	128	165527	29.938	ng/ul	0.00
24) 2-Nitrophenol-d4	9.811	143	186366	28.851	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.346	165	358537	26.689	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	368337	20.901	ng/ul	0.00
46) Dimethylphthalate-d6	13.965	166	1485354	30.709	ng/ul	0.00
49) Acenaphthylene-d8	14.253	160	1673854	30.851	ng/ul	0.00
54) 4-Nitrophenol-d4	14.764	143	39042m	5.636	ng/ul	0.00
60) Fluorene-d10	15.552	176	1342311	31.672	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	249213	27.369	ng/ul	0.00
73) Anthracene-d10	17.408	188	2379236	34.132	ng/ul	0.00
81) Pyrene-d10	19.700	212	2959404	34.757	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.518	264	2928226	35.104	ng/ul	0.00

Target Compounds Qvalue

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

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