

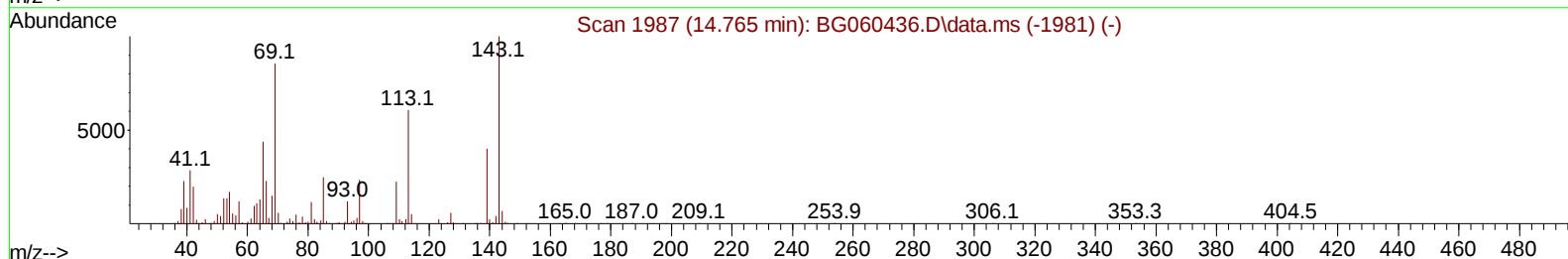
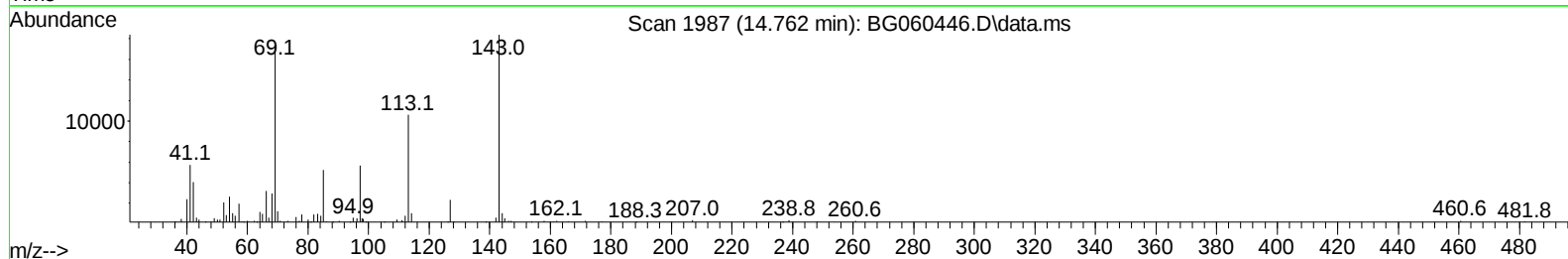
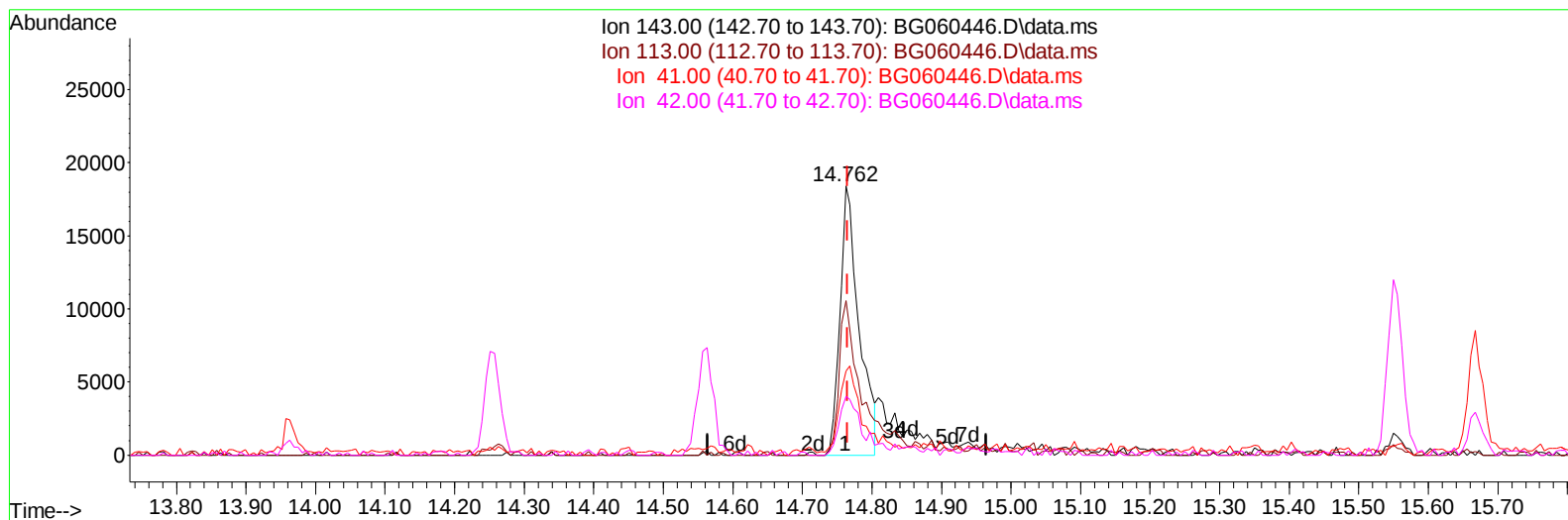
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012624\
 Data File : BG060446.D
 Acq On : 26 Jan 2024 18:18
 Operator : MA/JU
 Sample : P1238-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BH704

Manual IntegrationsAPPROVED

Quant Time: Jan 26 22:04:46 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: BG060446.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.762min (-0.002) 4.99 ng/u1

response 35018

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	57.50
41.00	29.80	31.23
42.00	22.70	21.93

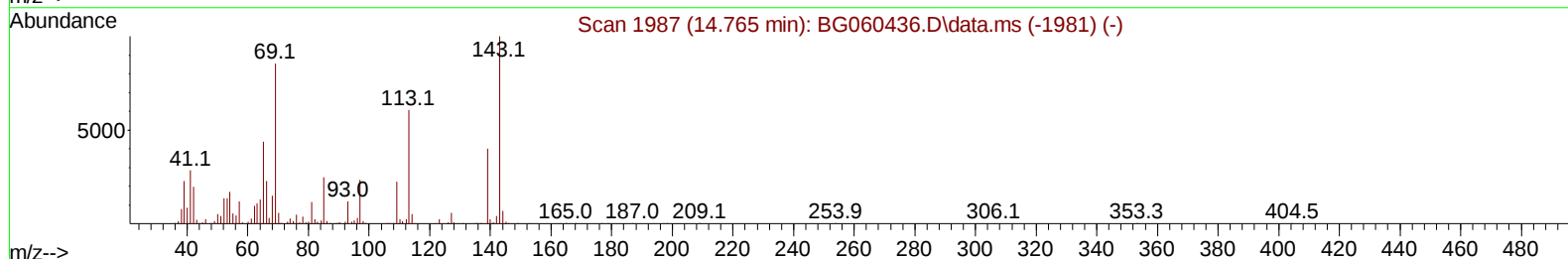
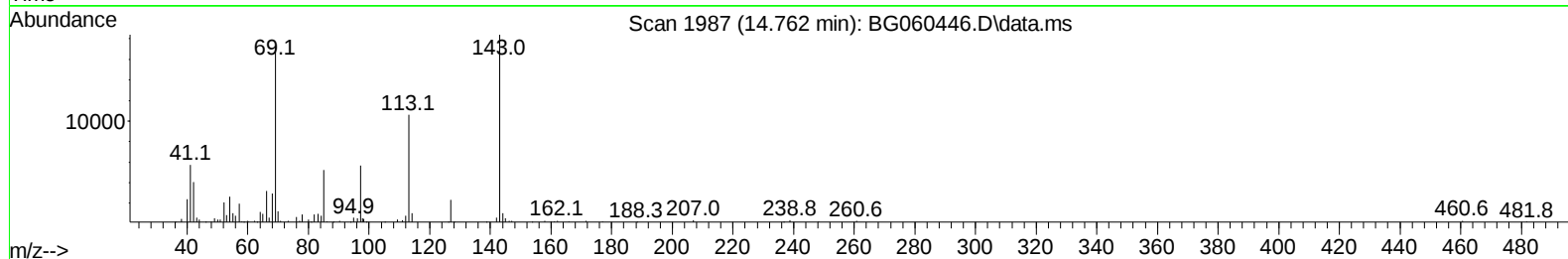
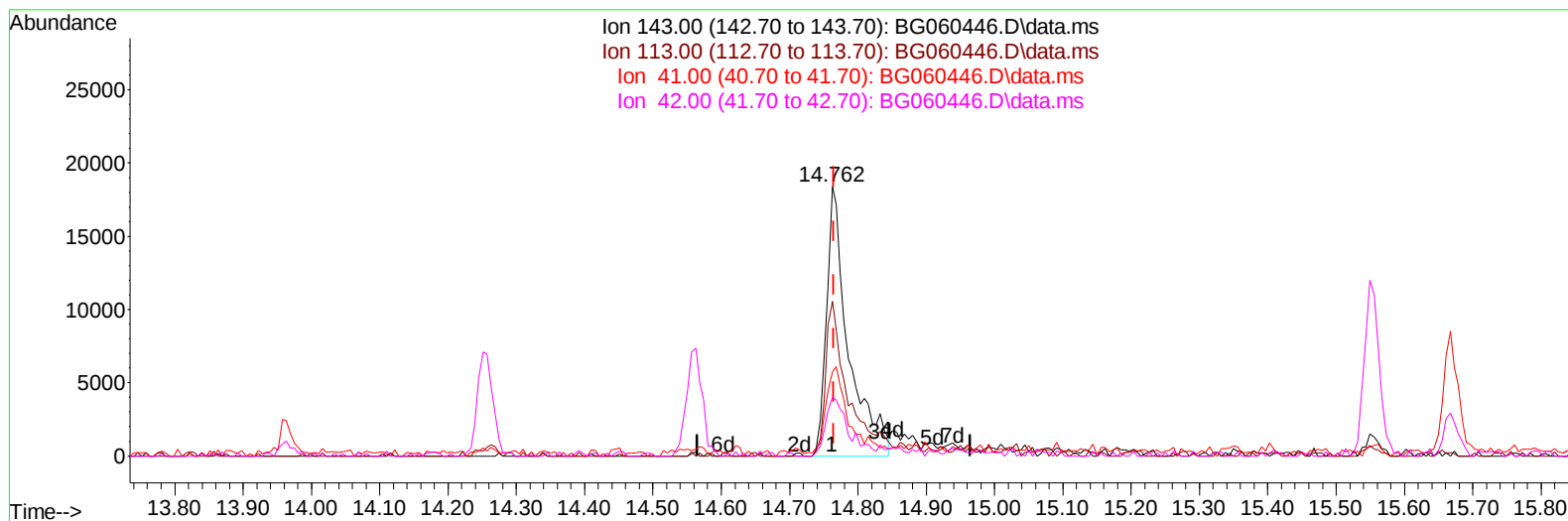
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TIC: BG060446.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.762min (-0.002) 5.87 ng/ul m

response 41226

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	57.50
41.00	29.80	31.23
42.00	22.70	21.93

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/29/2024
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Quant Time: Jan 26 22:13:49 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.923	152	170948	20.000	ng/ul	0.00
20) Naphthalene-d8	10.726	136	741735	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.563	164	599051	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.306	188	1478664	20.000	ng/ul	0.00
79) Chrysene-d12	21.572	240	1459129	20.000	ng/ul	0.00
88) Perylene-d12	24.733	264	1614112	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	20353	5.966	ng/uL	0.00
4) Pyridine-d5	3.775	84	128494	11.907	ng/ul	0.00
7) Phenol-d5	7.089	99	105695	6.510	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.253	67	252442	24.744	ng/ul	0.00
11) 2-Chlorophenol-d4	7.459	132	218527	20.156	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	153628	11.488	ng/ul	0.00
21) Nitrobenzene-d5	9.087	128	147814	27.015	ng/ul	0.00
24) 2-Nitrophenol-d4	9.809	143	172788	27.029	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.344	165	340065	25.579	ng/ul	0.00
31) 4-Chloroaniline-d4	10.867	131	381888	21.897	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166	1384751	28.249	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1517416	27.597	ng/ul	0.00
54) 4-Nitrophenol-d4	14.762	143	41226m	5.872	ng/ul	0.00
60) Fluorene-d10	15.555	176	1242642	28.931	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.667	200	228065	24.909	ng/ul	0.00
73) Anthracene-d10	17.406	188	2179459	31.093	ng/ul	0.00
81) Pyrene-d10	19.704	212	2771216	31.688	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.516	264	2727583	32.213	ng/ul	0.00

Target Compounds Qvalue

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

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