

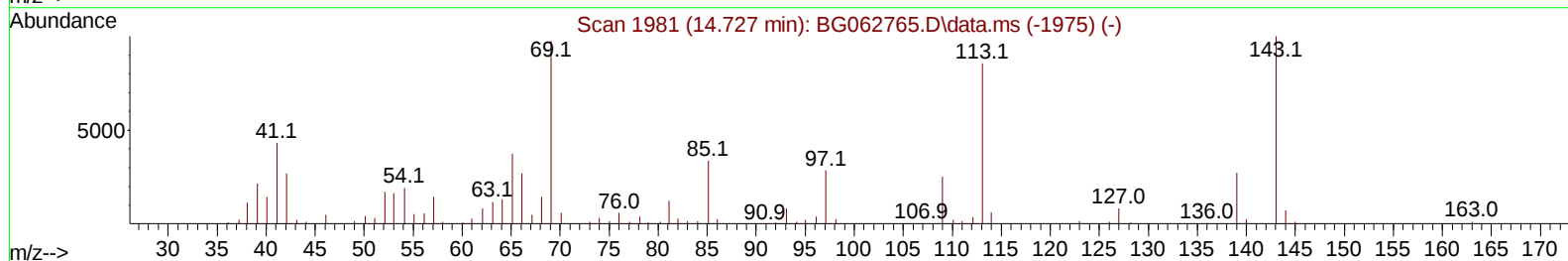
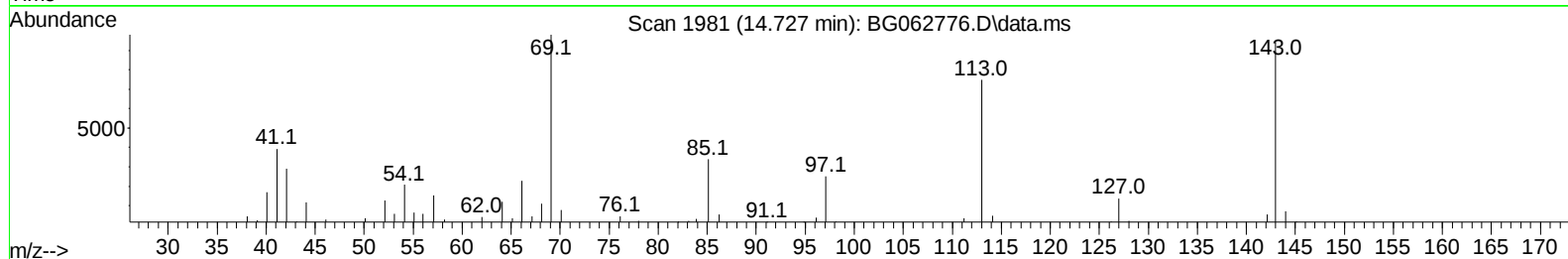
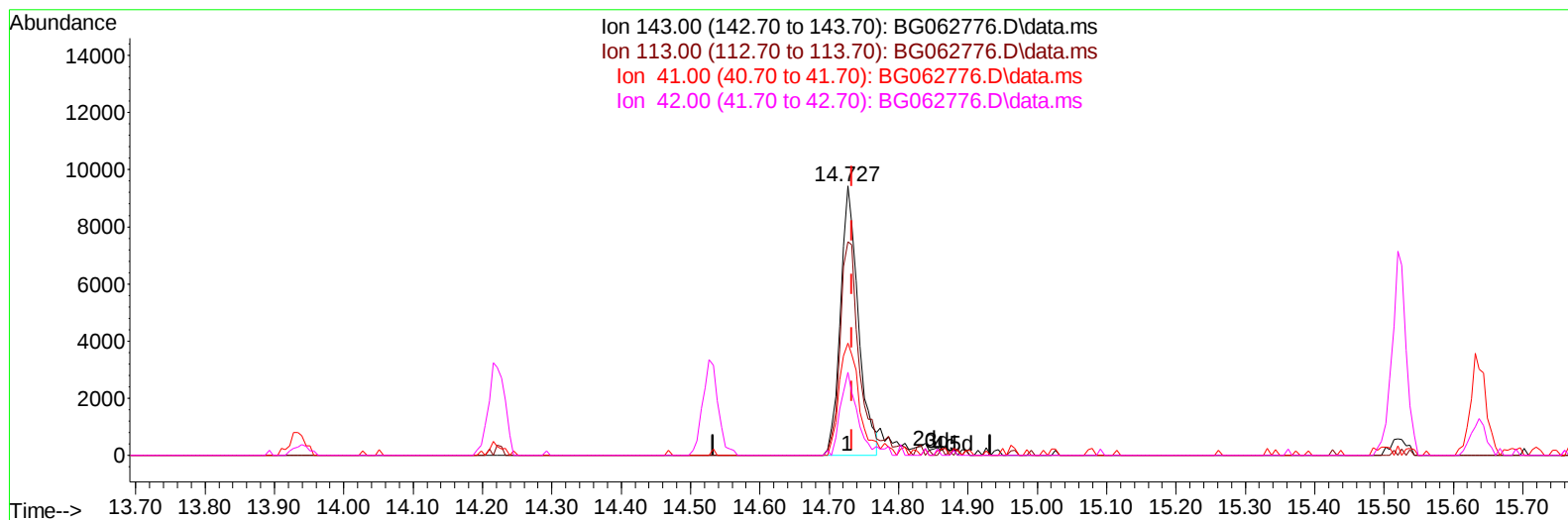
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062776.D
Acq On : 13 Sep 2024 00:15
Operator : JU/RC
Sample : P3922-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AW4

Manual IntegrationsAPPROVED

Quant Time: Sep 13 02:12:52 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 10 01:28:17 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/13/2024
Supervised By :mohammad ahmed 09/14/2024



TIC: BG062776.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.727min (-0.006) 7.57 ng/u1

response 16854

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.40	79.24
41.00	45.80	41.72
42.00	31.60	30.86

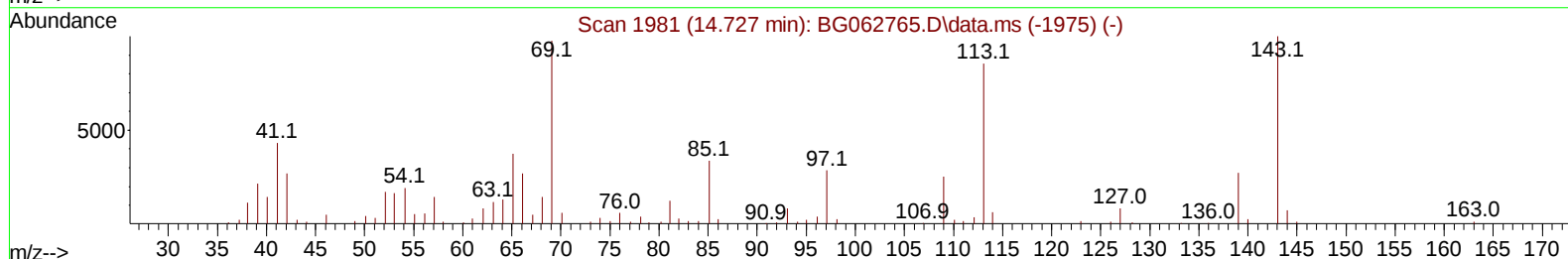
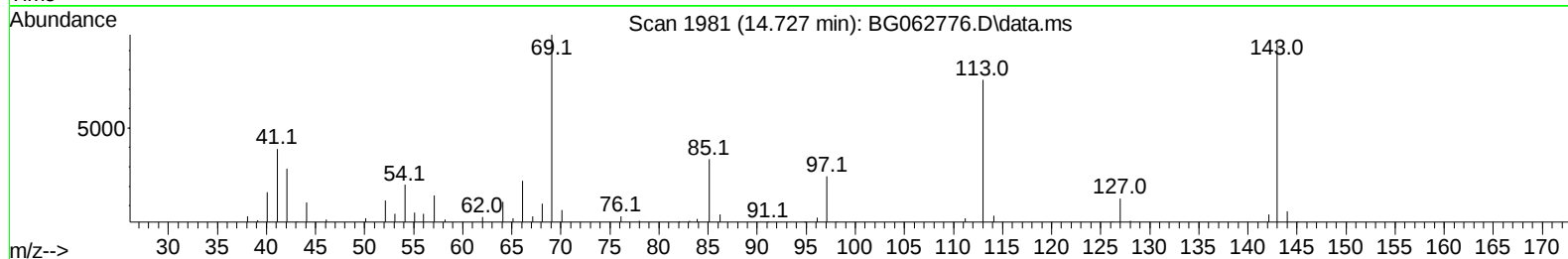
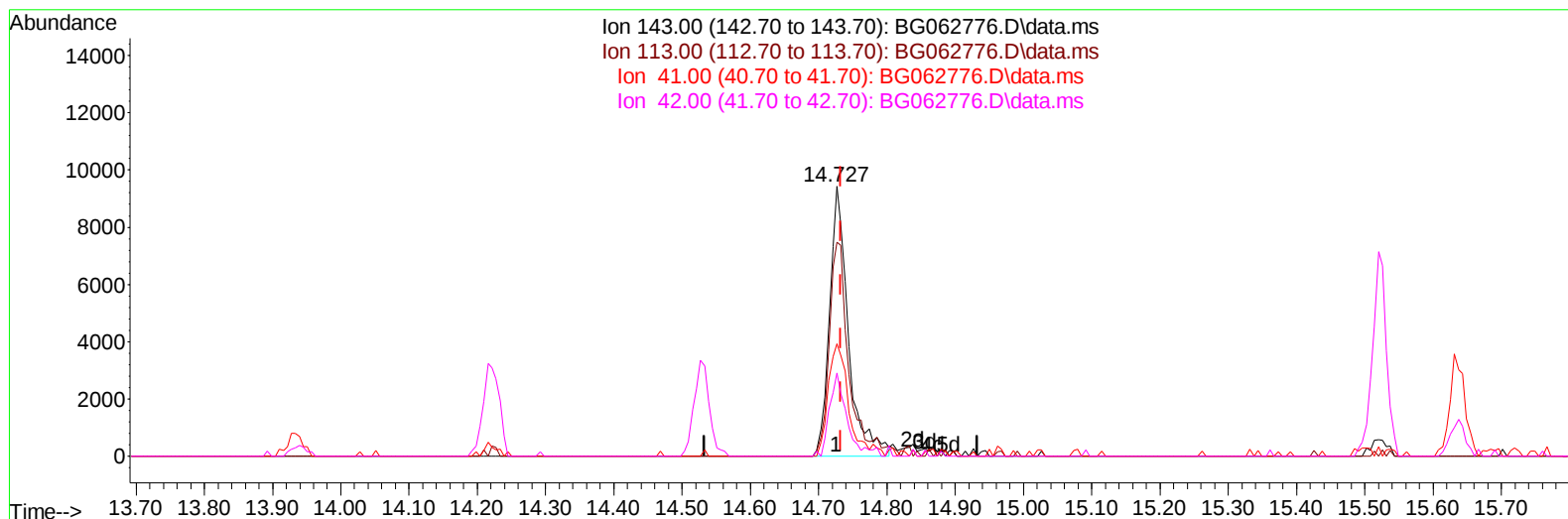
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TIC: BG062776.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.727min (-0.006) 8.09 ng/ul m

response 18005

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.40	79.24
41.00	45.80	41.72
42.00	31.60	30.86

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	43509	20.000	ng/ul	0.00
20) Naphthalene-d8	10.696	136	187512	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.533	164	170149	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.277	188	453904	20.000	ng/ul	0.00
79) Chrysene-d12	21.519	240	495072	20.000	ng/ul	0.00
88) Perylene-d12	24.580	264	543719	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	6066	6.302	ng/uL	0.00
4) Pyridine-d5	3.769	84	19963	6.516	ng/ul	0.00
7) Phenol-d5	7.065	99	37115	9.728	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.230	67	70251	30.712	ng/ul	0.00
11) 2-Chlorophenol-d4	7.435	132	81741	29.414	ng/ul	0.00
15) 4-Methylphenol-d8	8.604	113	70073	22.040	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	46109	34.405	ng/ul	0.00
24) 2-Nitrophenol-d4	9.774	143	55264	37.705	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.314	165	106230	33.281	ng/ul	0.00
31) 4-Chloroaniline-d4	10.825	131	92917	21.178	ng/ul	0.00
46) Dimethylphthalate-d6	13.933	166	424236	32.378	ng/ul	0.00
49) Acenaphthylene-d8	14.221	160	480741	33.036	ng/ul	0.00
54) 4-Nitrophenol-d4	14.727	143	18005m	8.089	ng/ul	0.00
60) Fluorene-d10	15.520	176	426058	36.085	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.637	200	82839	31.728	ng/ul	0.00
73) Anthracene-d10	17.371	188	825637	38.542	ng/ul	0.00
81) Pyrene-d10	19.662	212	1086862	39.013	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.374	264	1164760	40.570	ng/ul	0.00

Target Compounds Qvalue

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

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