

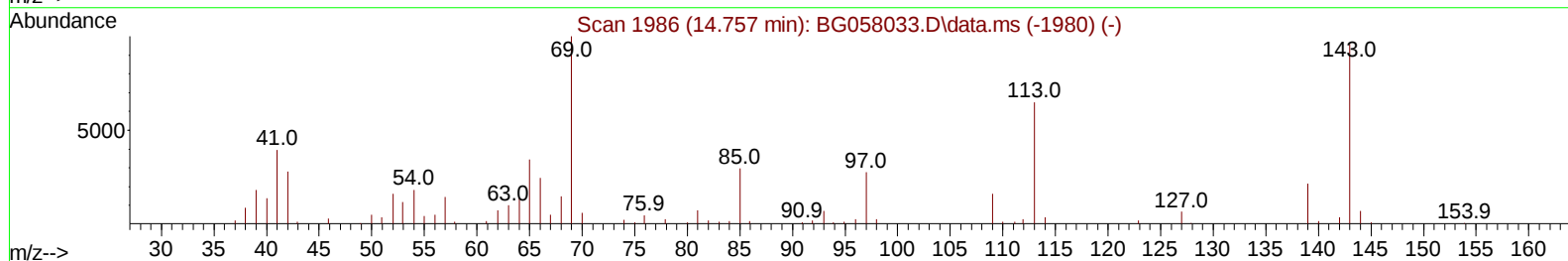
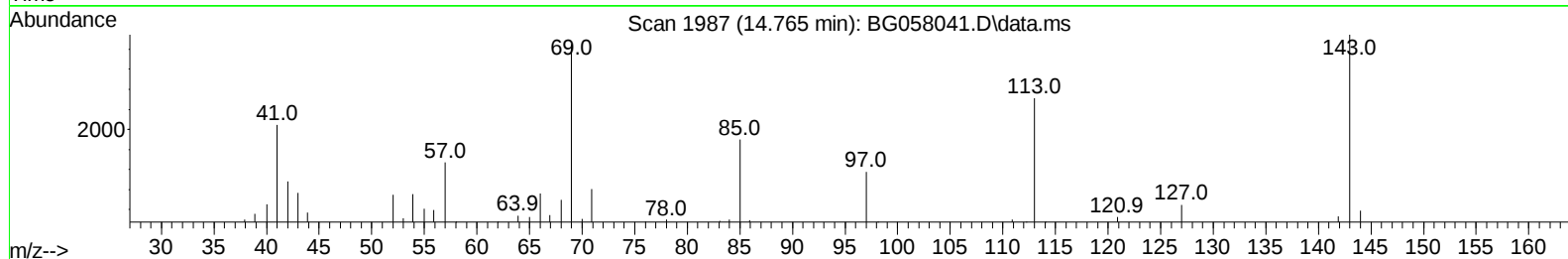
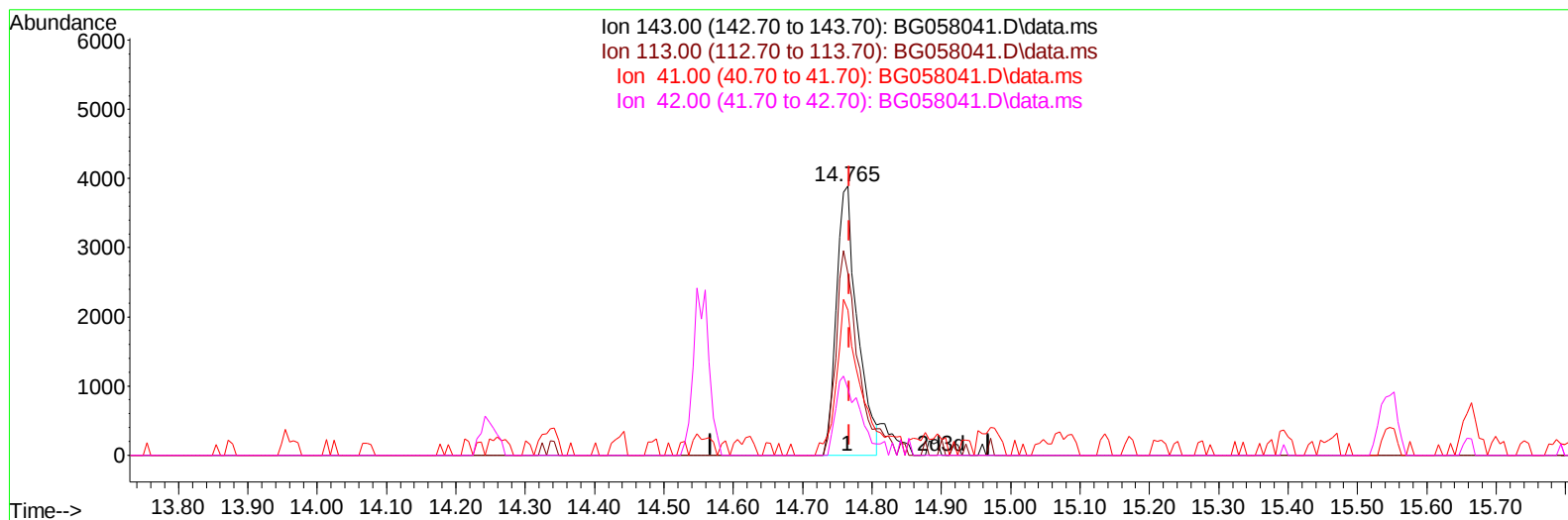
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070523\
Data File : BG058041.D
Acq On : 6 Jul 2023 8:38
Operator : CG/JU
Sample : 03244-20
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCGF9

Manual IntegrationsAPPROVED

Quant Time: Jul 06 09:16:41 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jul 05 23:20:13 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/06/2023
Supervised By :Sohil Jodhani 07/06/2023



TIC: BG058041.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.765min (-0.003) 4.60 ng/u1

response 8190

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.40	67.51
41.00	40.50	53.86#
42.00	27.50	24.68

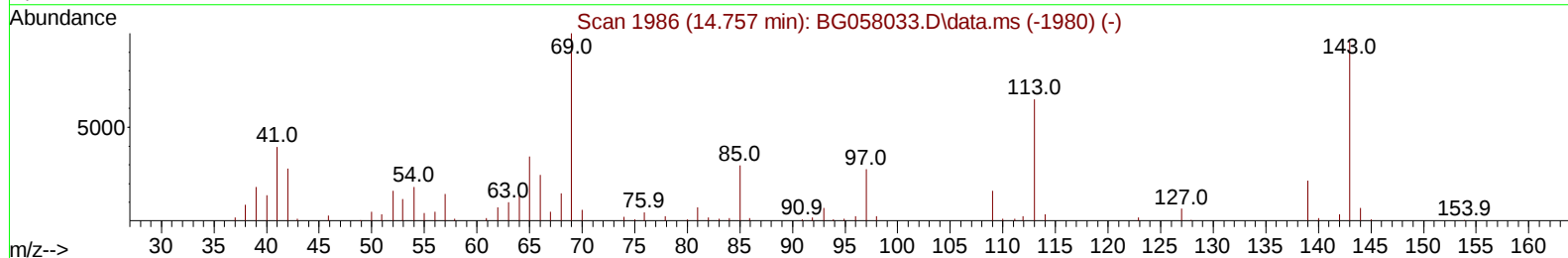
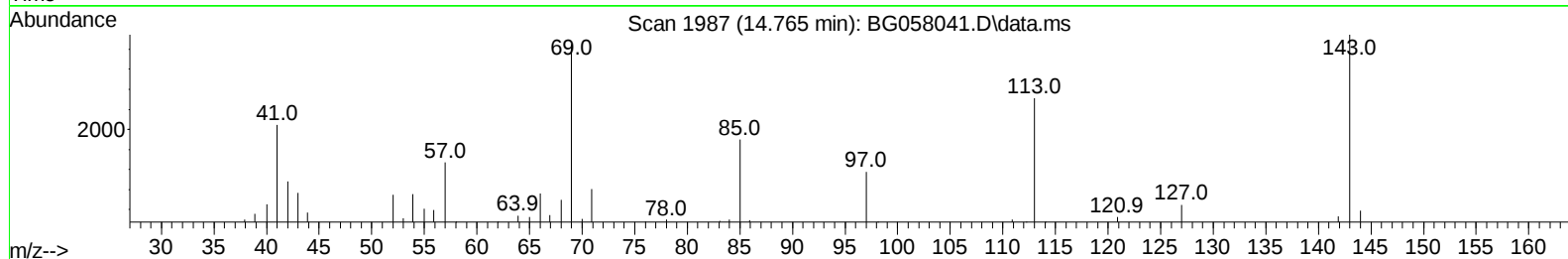
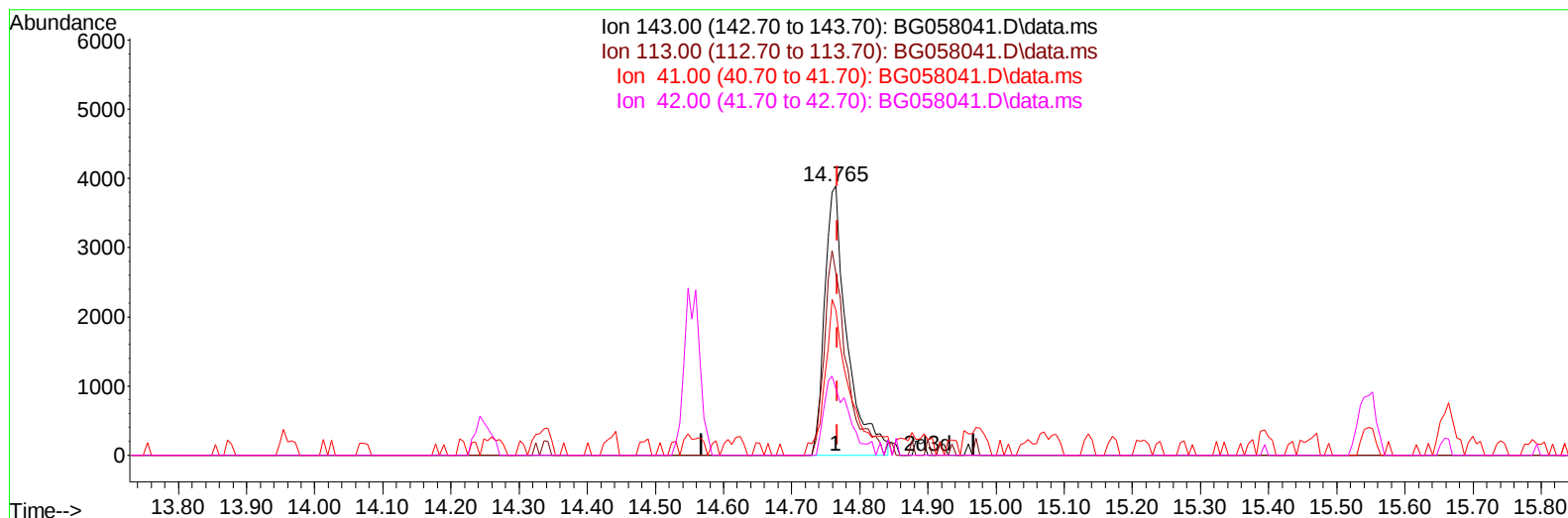
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TIC: BG058041.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.765min (-0.003) 4.98 ng/ul m

response 8869

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.40	67.51
41.00	40.50	53.86#
42.00	27.50	24.68

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/06/2023
 Supervised By :Sohil Jodhani 07/06/2023

Quant Time: Jul 06 09:18:46 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 05 23:20:13 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.920	152	37989	20.000	ng/ul	0.00
20) Naphthalene-d8	10.723	136	169062	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.554	164	127679	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.297	188	334198	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	312817	20.000	ng/ul	0.00
88) Perylene-d12	24.706	264	367808	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.361	96	1321	1.218	ng/uL	0.00
4) Pyridine-d5	3.772	84	3461	1.051	ng/ul	0.00
7) Phenol-d5	7.080	99	26112	6.731	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.244	67	17683	7.563	ng/ul	0.00
11) 2-Chlorophenol-d4	7.450	132	19499	7.287	ng/ul	0.00
15) 4-Methylphenol-d8	8.625	113	17239	5.889	ng/ul	0.00
21) Nitrobenzene-d5	9.078	128	10421	7.563	ng/ul	0.00
24) 2-Nitrophenol-d4	9.800	143	10801	7.458	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.341	165	19518	7.026	ng/ul	0.00
31) 4-Chloroaniline-d4	10.858	131	14274	3.389	ng/ul	0.00
46) Dimethylphthalate-d6	13.954	166	75306	7.475	ng/ul	0.00
49) Acenaphthylene-d8	14.248	160	82074	7.466	ng/ul	0.00
54) 4-Nitrophenol-d4	14.765	143	8869m	4.983	ng/ul	0.00
60) Fluorene-d10	15.546	176	69330	8.104	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.664	200	8042	4.432	ng/ul	0.00
73) Anthracene-d10	17.397	188	100987	6.512	ng/ul	0.00
81) Pyrene-d10	19.689	212	158896	8.065	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.489	264	145129	7.745	ng/ul	0.00

Target Compounds Qvalue

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

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