

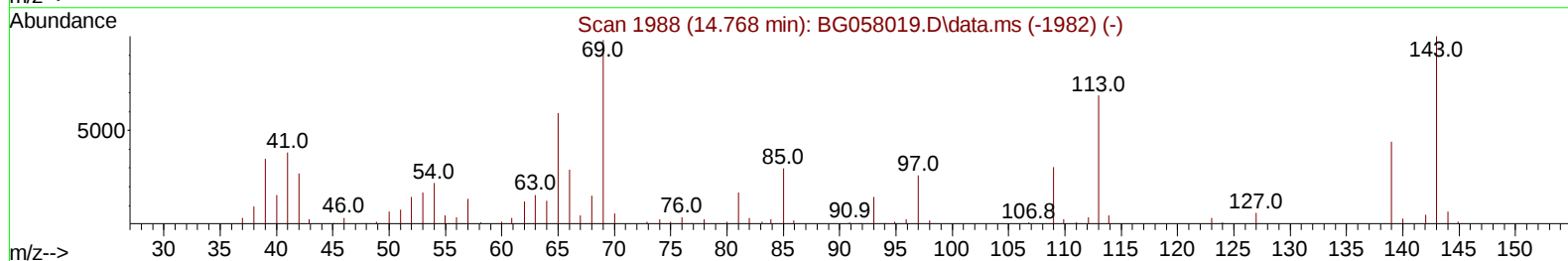
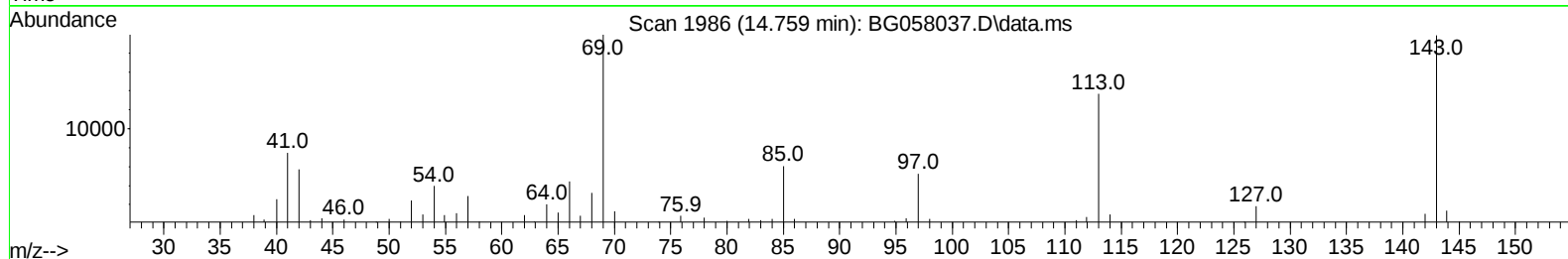
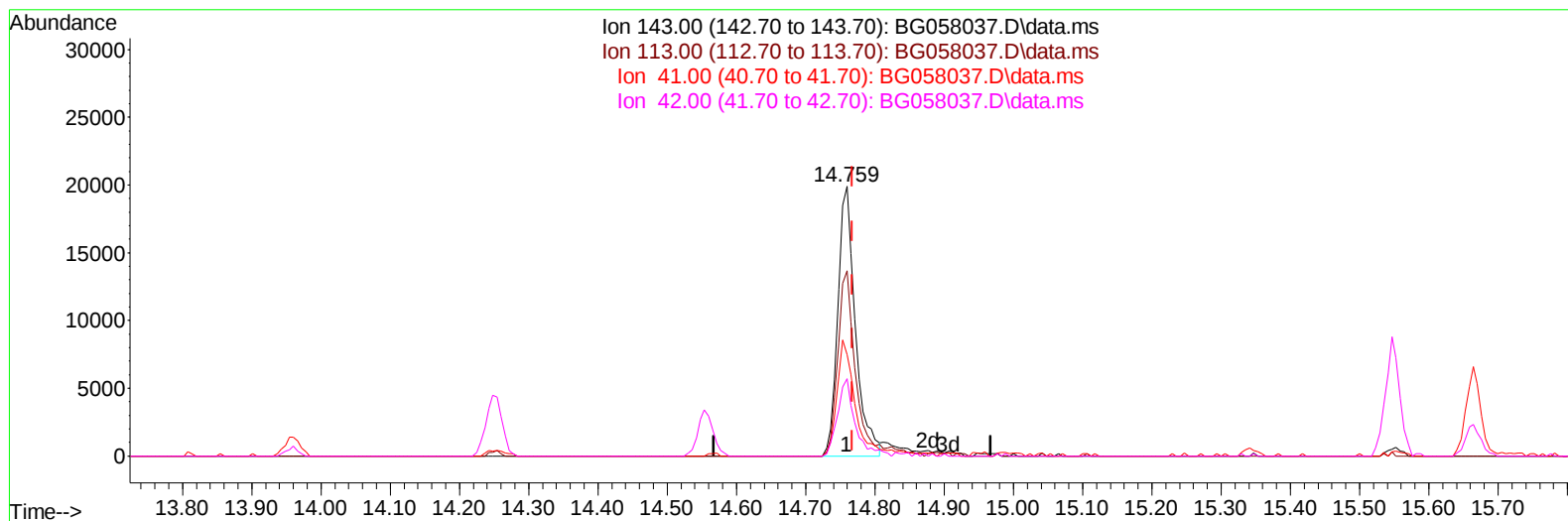
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070523\
Data File : BG058037.D
Acq On : 6 Jul 2023 5:50
Operator : CG/JU
Sample : 03257-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EXYY3

Manual IntegrationsAPPROVED

Quant Time: Jul 06 06:24:38 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: BG058037.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.759min (-0.008) 13.58 ng/u1

response 34977

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.40	68.70
41.00	40.50	37.64
42.00	27.50	28.70

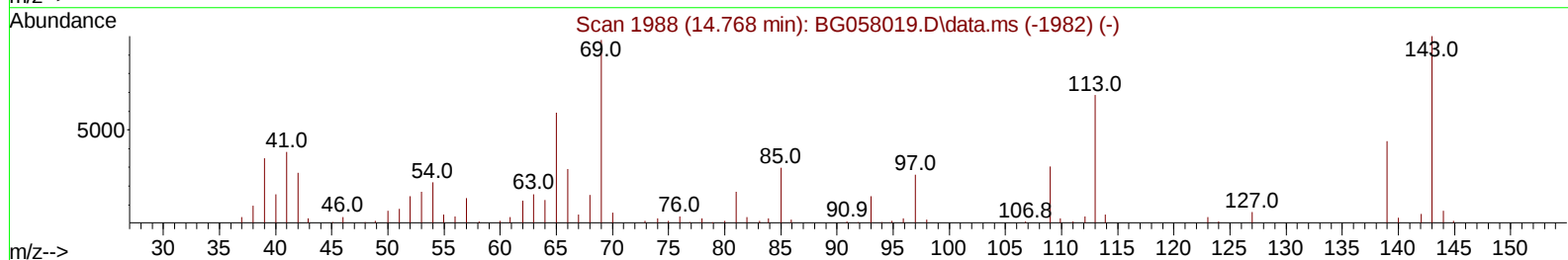
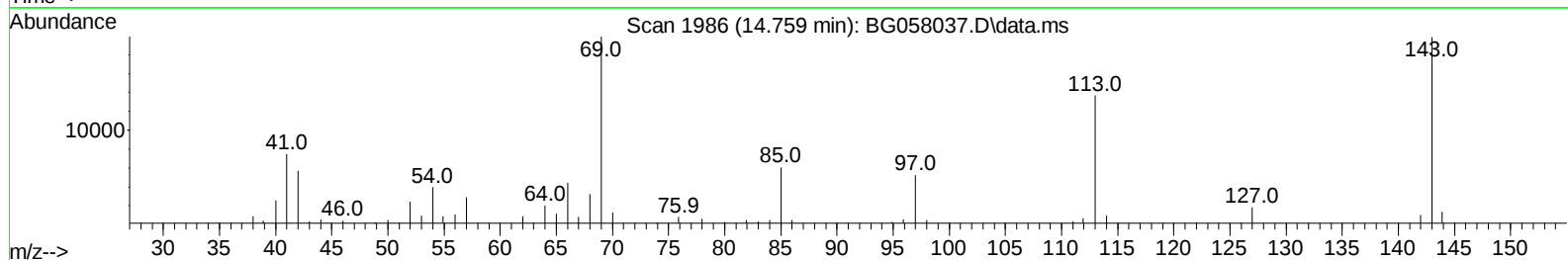
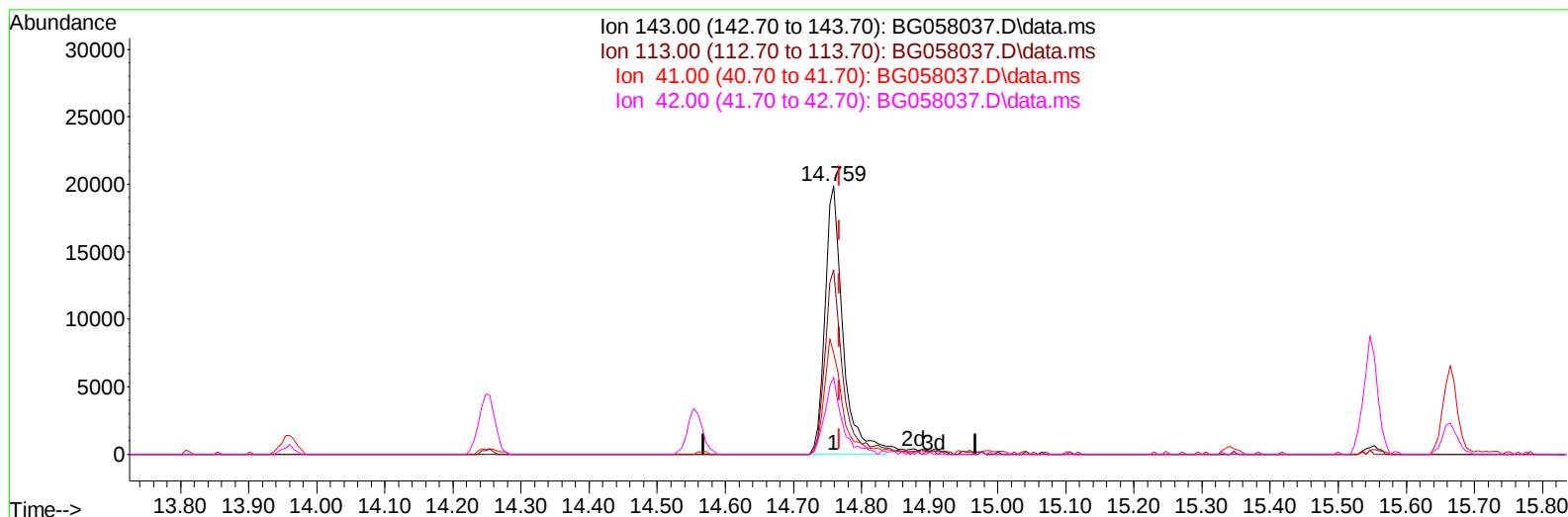
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070523\
Data File : BG058037.D
Acq On : 6 Jul 2023 5:50
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Sample : 03257-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EXYY3

Manual IntegrationsAPPROVED

Quant Time: Jul 06 06:24:38 2023
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TIC: BG058037.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.759min (-0.008) 14.13 ng/ul m

response 36402

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.40	68.70
41.00	40.50	37.64
42.00	27.50	28.70

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070523\
 Data File : BG058037.D
 Acq On : 6 Jul 2023 5:50
 Operator : CG/JU
 Sample : 03257-07
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EXYY3

Manual IntegrationsAPPROVED

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

Quant Time: Jul 06 06:25:13 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	59487	20.000	ng/ul	0.00
20) Naphthalene-d8	10.723	136	259925	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.554	164	184822	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.303	188	481556	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	458015	20.000	ng/ul	0.00
88) Perylene-d12	24.712	264	577891	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.355	96	9983	5.876	ng/uL	0.00
4) Pyridine-d5	3.766	84	76725	14.877	ng/ul	0.00
7) Phenol-d5	7.086	99	74131	12.202	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.245	67	133545	36.477	ng/ul	0.00
11) 2-Chlorophenol-d4	7.450	132	121060	28.893	ng/ul	0.00
15) 4-Methylphenol-d8	8.625	113	93745	20.450	ng/ul	0.00
21) Nitrobenzene-d5	9.078	128	82485	38.936	ng/ul	0.00
24) 2-Nitrophenol-d4	9.800	143	87187	39.158	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.341	165	142552	33.375	ng/ul	0.00
31) 4-Chloroaniline-d4	10.858	131	212023	32.741	ng/ul	0.00
46) Dimethylphthalate-d6	13.960	166	630747	43.253	ng/ul	0.00
49) Acenaphthylene-d8	14.248	160	639711	40.200	ng/ul	0.00
54) 4-Nitrophenol-d4	14.759	143	36402m	14.129	ng/ul	0.00
60) Fluorene-d10	15.547	176	538407	43.474	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.664	200	118278	45.233	ng/ul	0.00
73) Anthracene-d10	17.403	188	1007633	45.094	ng/ul	0.00
81) Pyrene-d10	19.689	212	1230213	42.648	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.501	264	1357786	46.117	ng/ul	0.00

Target Compounds Qvalue

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

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