

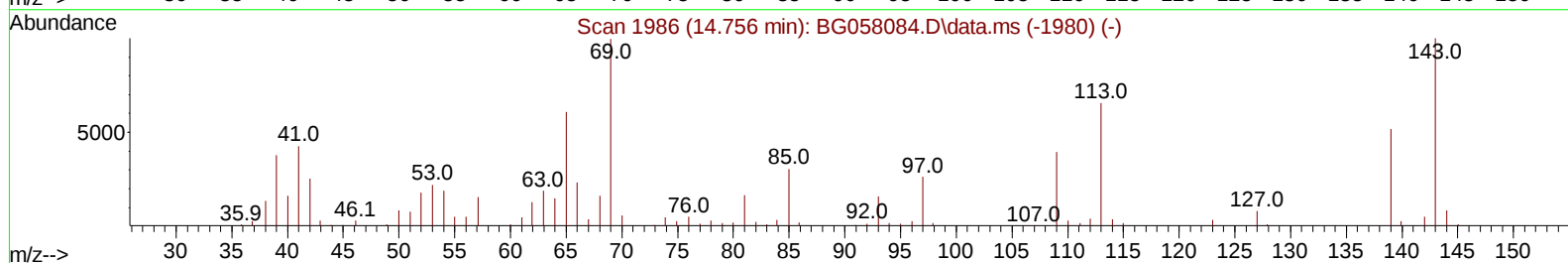
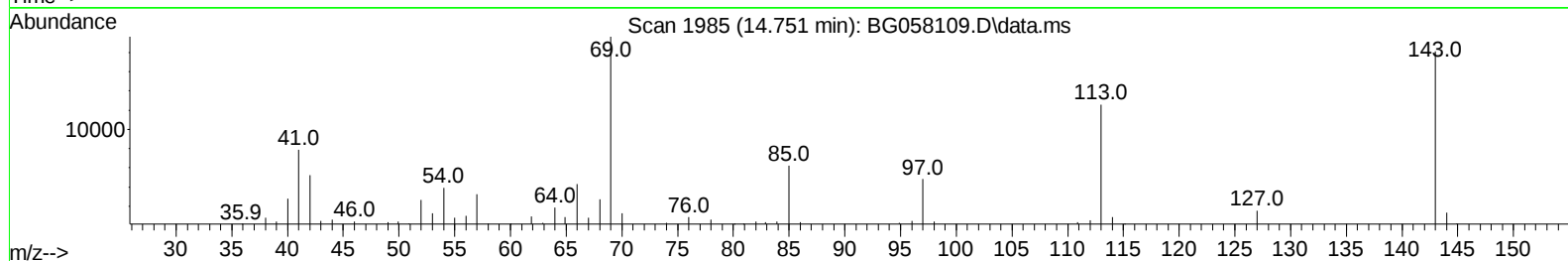
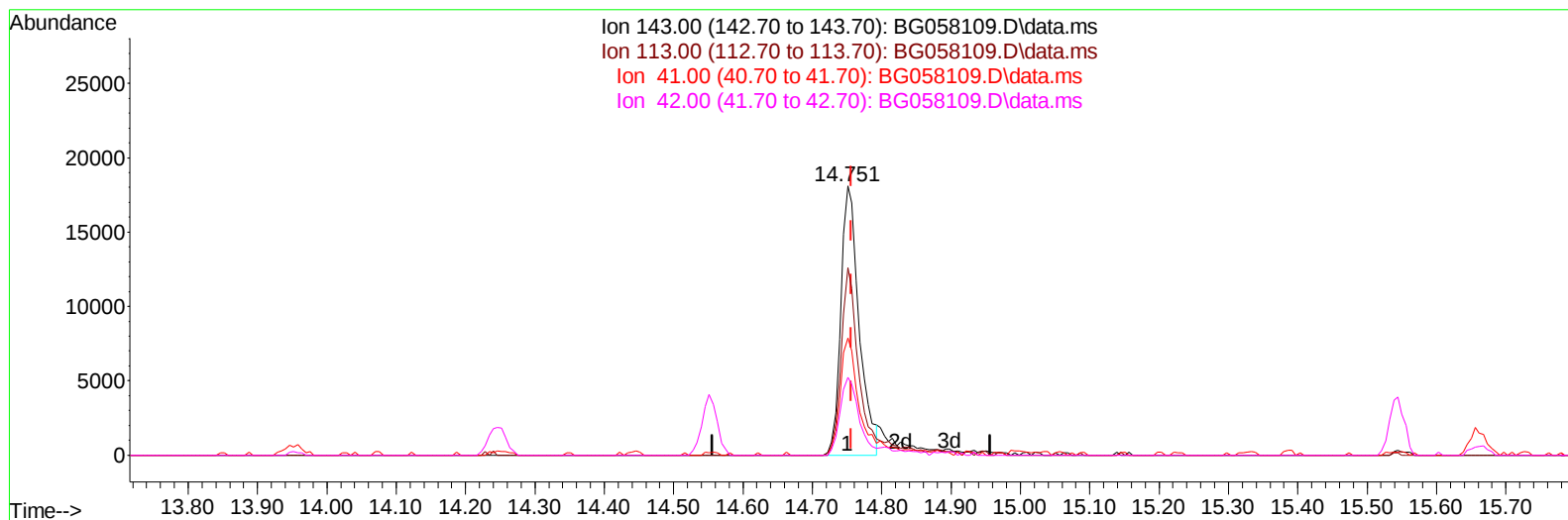
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070823\
 Data File : BG058109.D
 Acq On : 9 Jul 2023 2:14
 Operator : CG/JU
 Sample : 03349-02
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YBTZ6

Manual IntegrationsAPPROVED

Quant Time: Jul 09 02:50:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 08 10:01:01 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/09/2023
 Supervised By :mohammad ahmed 07/09/2023



TIC: BG058109.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (-0.005) 11.12 ng/u1

response 32953

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.40	69.57
41.00	40.50	43.41
42.00	27.50	28.89

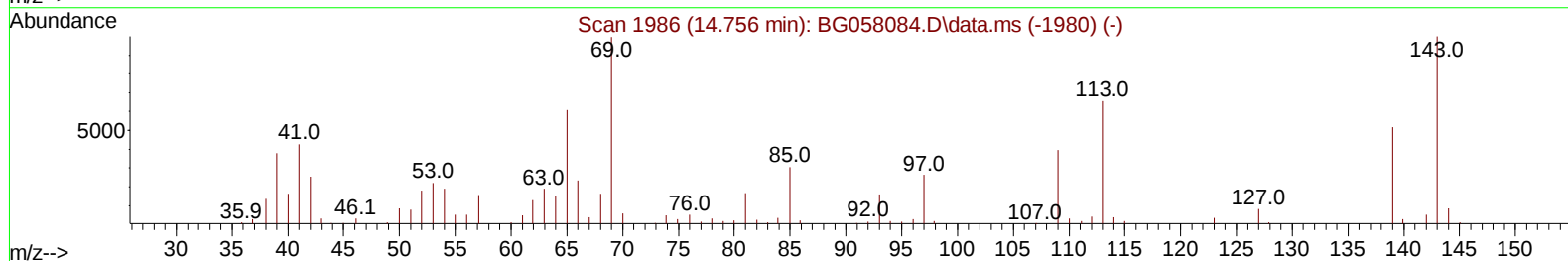
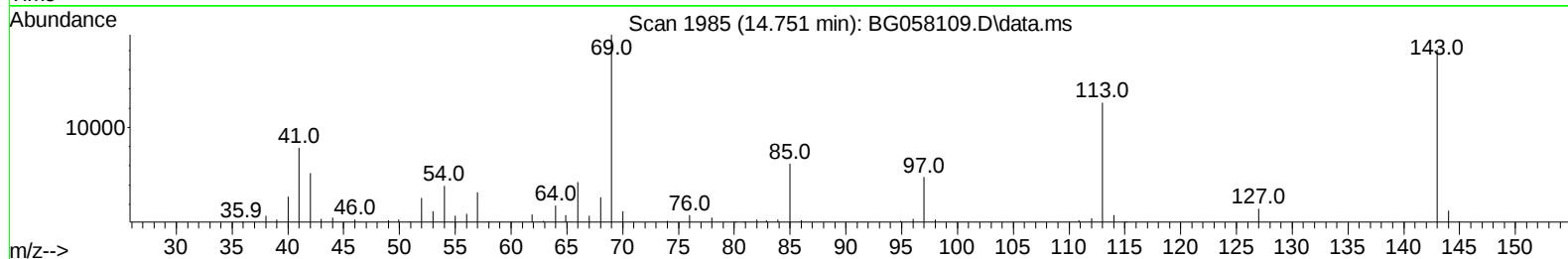
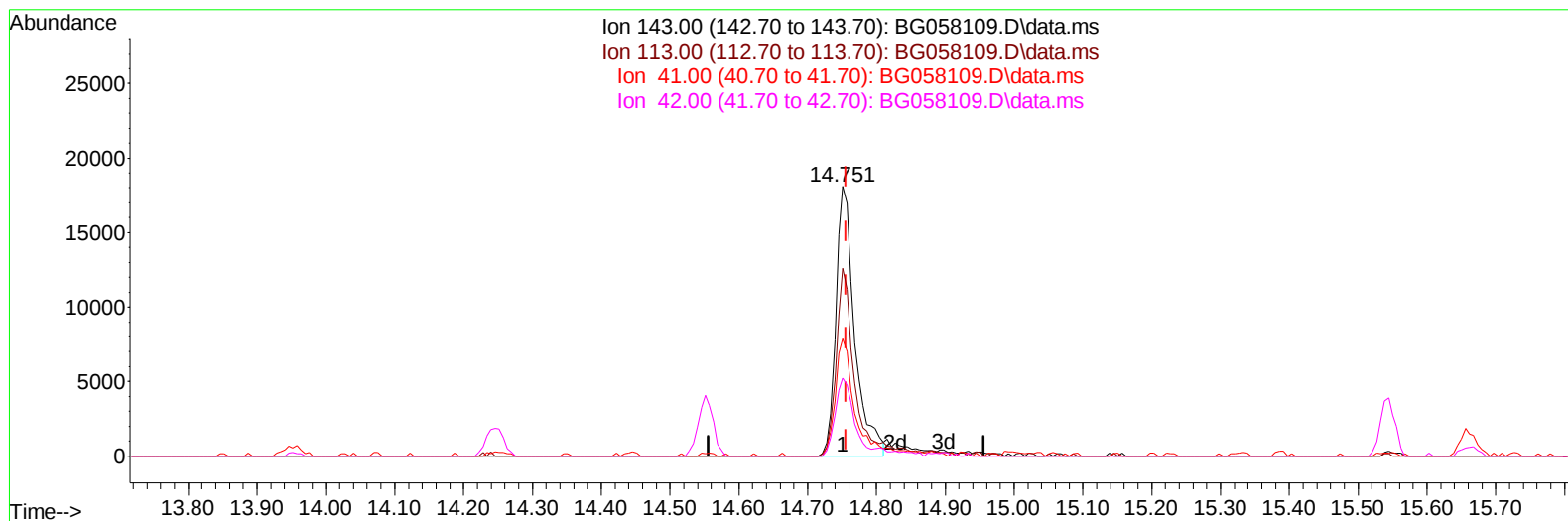
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TIC: BG058109.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (-0.005) 11.60 ng/ul m

response 34375

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.40	69.57
41.00	40.50	43.41
42.00	27.50	28.89

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.918	152	59667	20.000	ng/ul	0.00
20) Naphthalene-d8	10.721	136	286439	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	212639	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.301	188	520115	20.000	ng/ul	0.00
79) Chrysene-d12	21.549	240	463000	20.000	ng/ul	0.00
88) Perylene-d12	24.704	264	565710	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.359	96	5210	3.058	ng/uL	0.00
4) Pyridine-d5	3.764	84	64496	12.468	ng/ul	0.00
7) Phenol-d5	7.078	99	97469	15.996	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.242	67	61508	16.750	ng/ul	0.00
11) 2-Chlorophenol-d4	7.448	132	68017	16.184	ng/ul	0.00
15) 4-Methylphenol-d8	8.623	113	72511	15.771	ng/ul	0.00
21) Nitrobenzene-d5	9.075	128	37604	16.107	ng/ul	0.00
24) 2-Nitrophenol-d4	9.798	143	41371	16.861	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.339	165	76093	16.166	ng/ul	0.00
31) 4-Chloroaniline-d4	10.856	131	106709	14.953	ng/ul	0.00
46) Dimethylphthalate-d6	13.952	166	266982	15.913	ng/ul	0.00
49) Acenaphthylene-d8	14.246	160	308477	16.849	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	34375m	11.597	ng/ul	0.00
60) Fluorene-d10	15.544	176	241983	16.983	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.662	200	33647	11.914	ng/ul	0.00
73) Anthracene-d10	17.395	188	404598	16.764	ng/ul	0.00
81) Pyrene-d10	19.686	212	470340	16.130	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.487	264	479024	16.620	ng/ul	0.00

Target Compounds Qvalue

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

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