

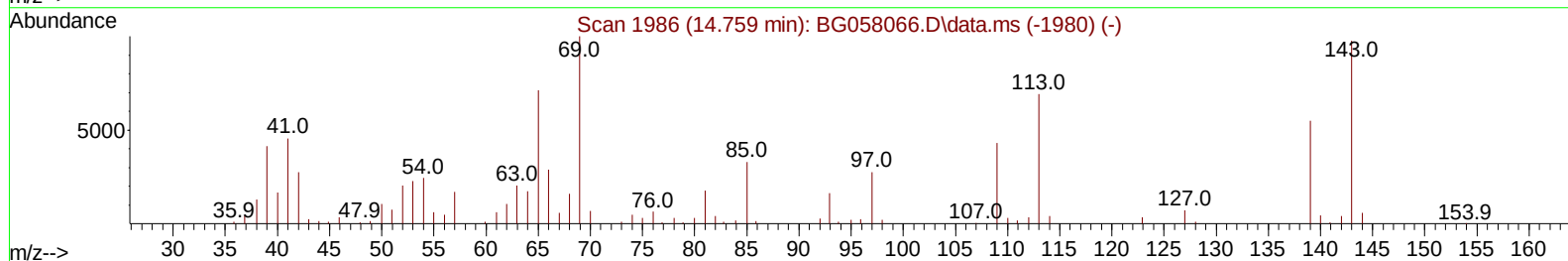
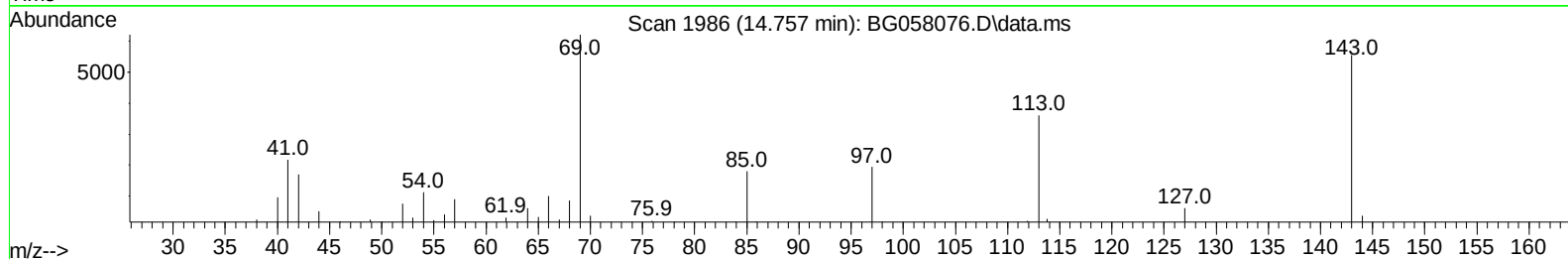
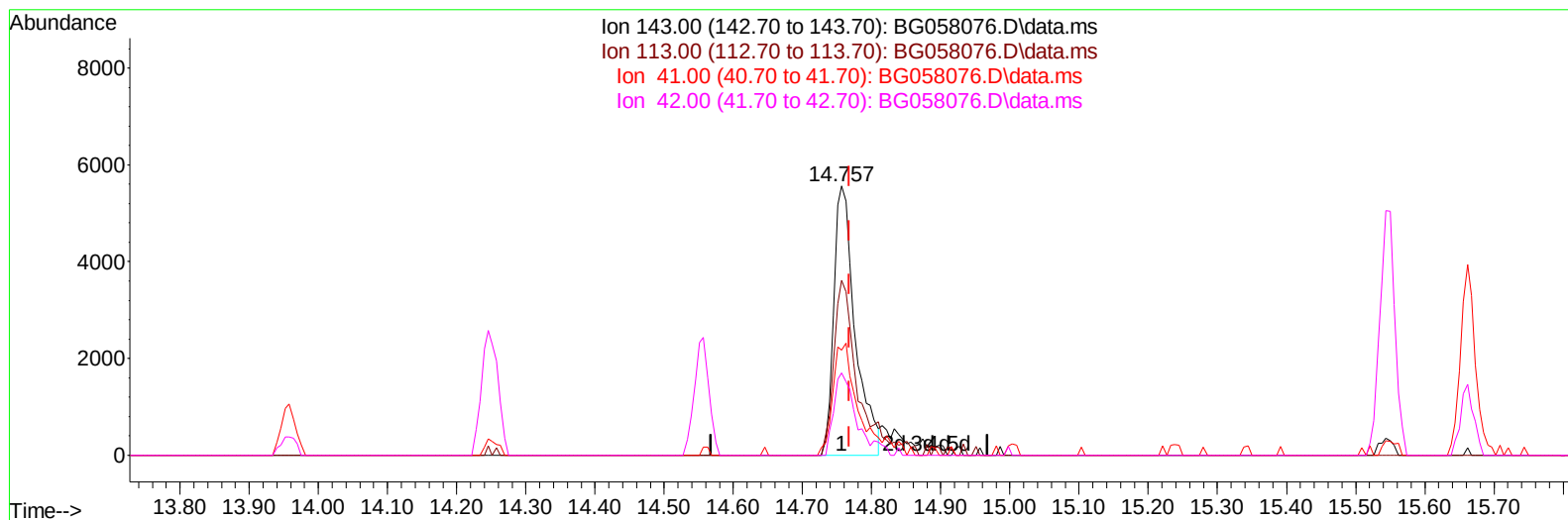
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070523\
Data File : BG058076.D
Acq On : 7 Jul 2023 15:31
Operator : CG/JU
Sample : 03284-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCGZ2

Manual IntegrationsAPPROVED

Quant Time: Jul 07 18:49:04 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 07/07/2023
Supervised By :Sohil Jodhani 07/07/2023



TIC: BG058076.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (-0.011) 6.51 ng/uL

response 11881

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.40	64.86
41.00	40.50	39.15
42.00	27.50	30.44

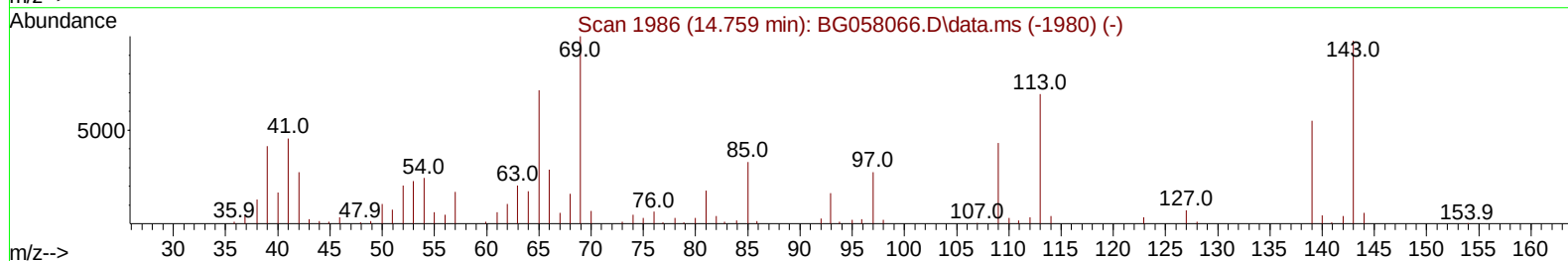
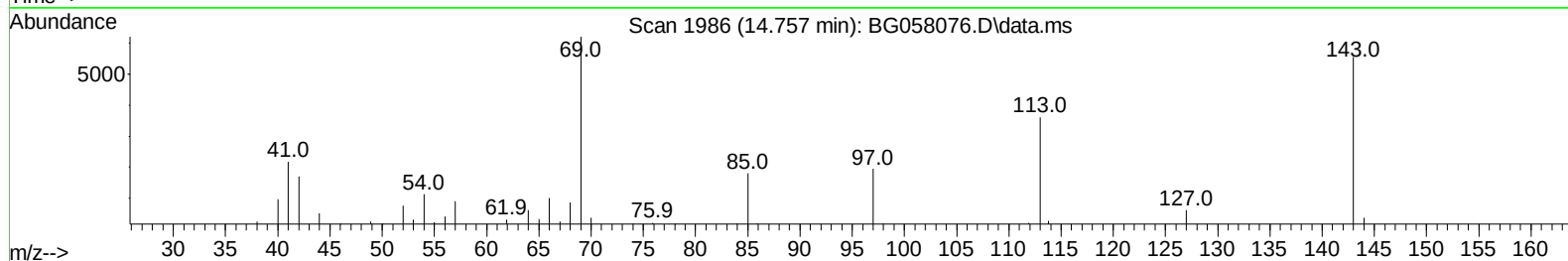
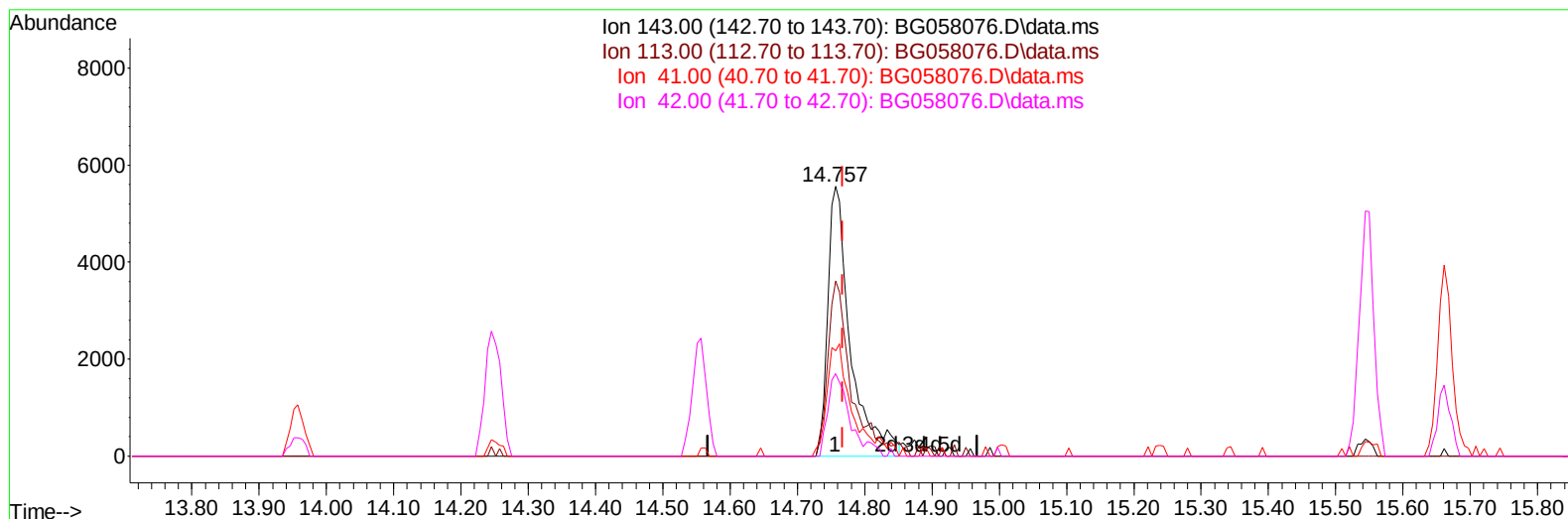
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(54) 4-Nitrophenol-d4 (S)

14.757min (-0.011) 7.09 ng/ul m

response 12939

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.40	64.86
41.00	40.50	39.15
42.00	27.50	30.44

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.918	152	37696	20.000	ng/ul	0.00
20) Naphthalene-d8	10.720	136	174167	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.557	164	130868	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.301	188	342404	20.000	ng/ul	0.00
79) Chrysene-d12	21.555	240	320587	20.000	ng/ul	0.00
88) Perylene-d12	24.710	264	392171	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	4805	4.463	ng/uL	0.00
4) Pyridine-d5	3.770	84	7331	2.243	ng/ul	0.00
7) Phenol-d5	7.077	99	27363	7.108	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.242	67	83658	36.060	ng/ul	0.00
11) 2-Chlorophenol-d4	7.448	132	71592	26.964	ng/ul	-0.01
15) 4-Methylphenol-d8	8.623	113	49981	17.206	ng/ul	-0.01
21) Nitrobenzene-d5	9.075	128	49189	34.651	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.804	143	53289	35.718	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.338	165	84629	29.570	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.855	131	54002	12.445	ng/ul	0.00
46) Dimethylphthalate-d6	13.958	166	405829	39.303	ng/ul	0.00
49) Acenaphthylene-d8	14.251	160	407354	36.152	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	12939m	7.093	ng/ul	-0.01
60) Fluorene-d10	15.550	176	342555	39.063	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.662	200	65820	35.401	ng/ul	-0.01
73) Anthracene-d10	17.401	188	645466	40.625	ng/ul	0.00
81) Pyrene-d10	19.692	212	810356	40.135	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.498	264	848575	42.471	ng/ul	0.00

Target Compounds Qvalue

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

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