

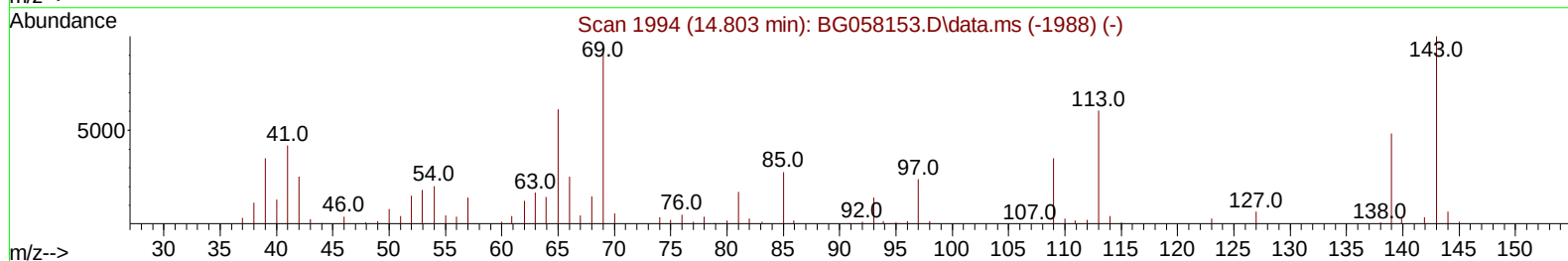
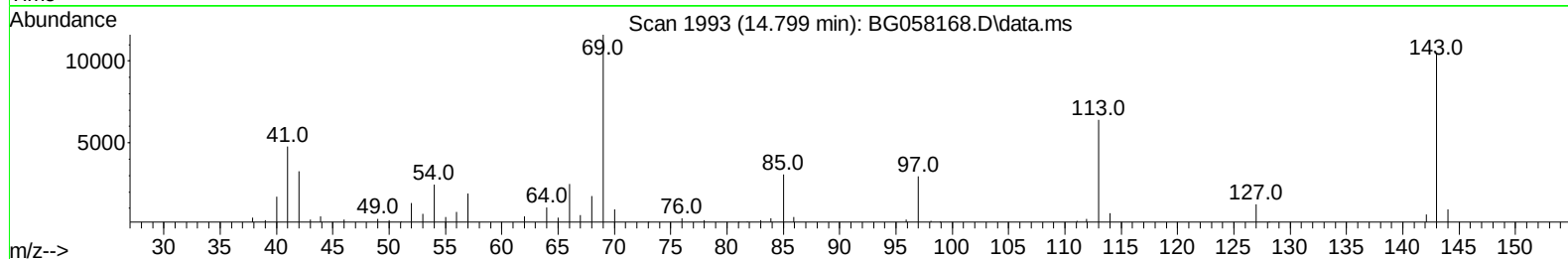
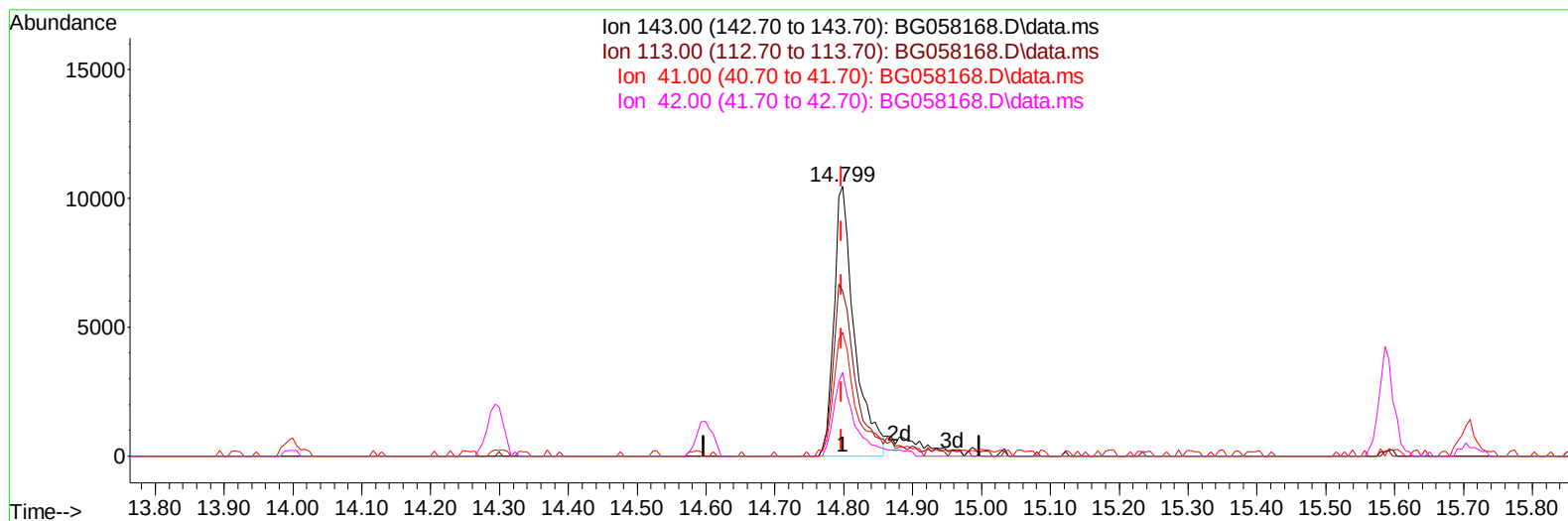
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071123\
Data File : BG058168.D
Acq On : 11 Jul 2023 21:10
Operator : CG/JU
Sample : 03385-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EX7Y5

Manual IntegrationsAPPROVED

Quant Time: Jul 11 22:58:20 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 10 16:52:40 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/12/2023
Supervised By :mohammad ahmed 07/12/2023



TIC: BG058168.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.799min (+ 0.001) 18.67 ng/ul

response 21729

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.40	61.31
41.00	40.50	45.83
42.00	27.50	31.09

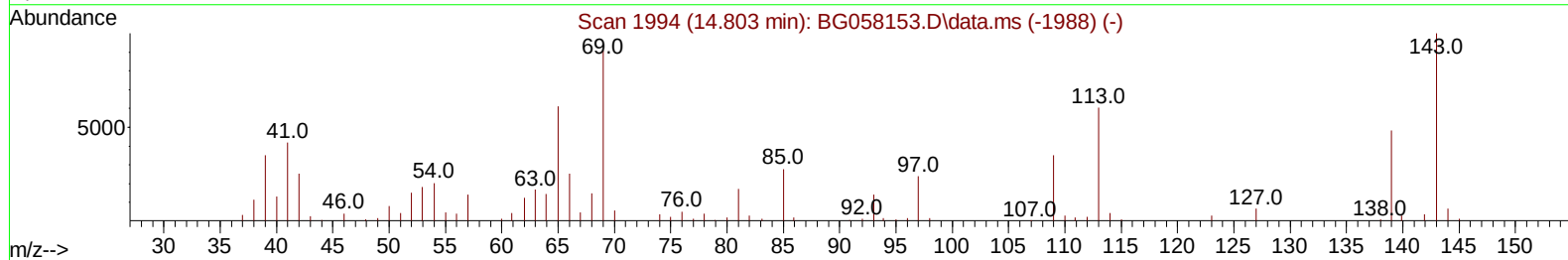
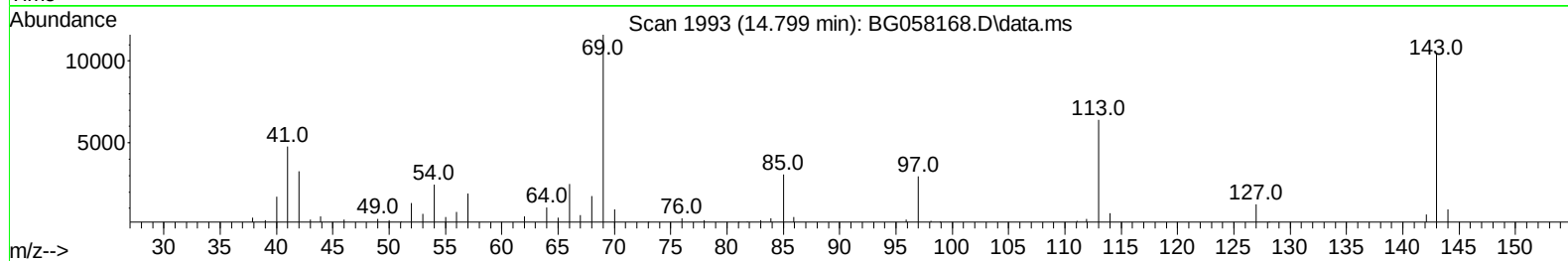
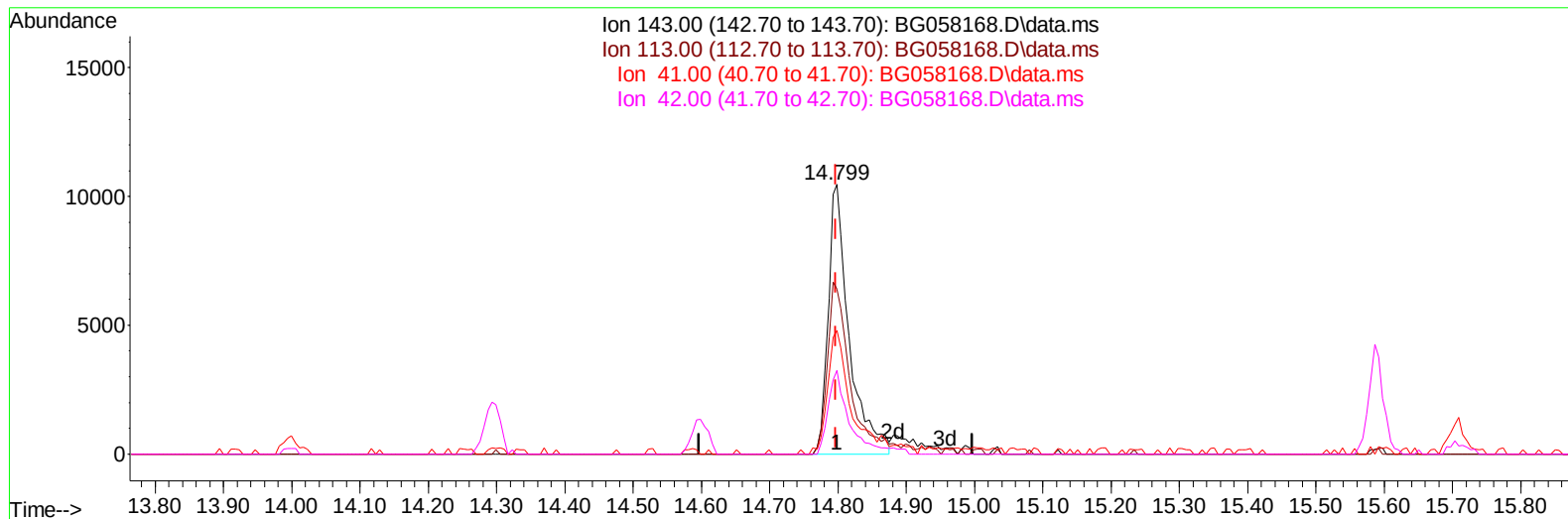
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TIC: BG058168.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.799min (+ 0.001) 19.26 ng/ul m

response 22426

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.40	61.31
41.00	40.50	45.83
42.00	27.50	31.09

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 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EX7Y5

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/12/2023
 Supervised By :mohammad ahmed 07/12/2023

Quant Time: Jul 11 23:43:11 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 10 16:52:40 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.971	152	25515	20.000	ng/ul	0.00
20) Naphthalene-d8	10.774	136	118215	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.599	164	83508	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.343	188	191649	20.000	ng/ul	0.00
79) Chrysene-d12	21.596	240	169128	20.000	ng/ul	0.00
88) Perylene-d12	24.793	264	193589	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.400	96	4956	6.802	ng/uL	0.00
4) Pyridine-d5	3.811	84	77902	35.216	ng/ul	0.00
7) Phenol-d5	7.131	99	100948	38.741	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.296	67	60129	38.292	ng/ul	0.00
11) 2-Chlorophenol-d4	7.507	132	69316	38.570	ng/ul	0.00
15) 4-Methylphenol-d8	8.676	113	78659	40.007	ng/ul	0.00
21) Nitrobenzene-d5	9.135	128	37788	39.219	ng/ul	0.00
24) 2-Nitrophenol-d4	9.857	143	43019	42.482	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	80931	41.662	ng/ul	0.00
31) 4-Chloroaniline-d4	10.909	131	108864	36.963	ng/ul	0.00
46) Dimethylphthalate-d6	13.999	166	277709	42.148	ng/ul	0.00
49) Acenaphthylene-d8	14.293	160	313403	43.589	ng/ul	0.00
54) 4-Nitrophenol-d4	14.799	143	22426m	19.265	ng/ul	0.00
60) Fluorene-d10	15.592	176	245141	43.809	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.709	200	20063	19.279	ng/ul	0.00
73) Anthracene-d10	17.442	188	391395	44.012	ng/ul	0.00
81) Pyrene-d10	19.728	212	452116	42.445	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.569	264	471160	47.771	ng/ul	0.00

Target Compounds Qvalue

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

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