

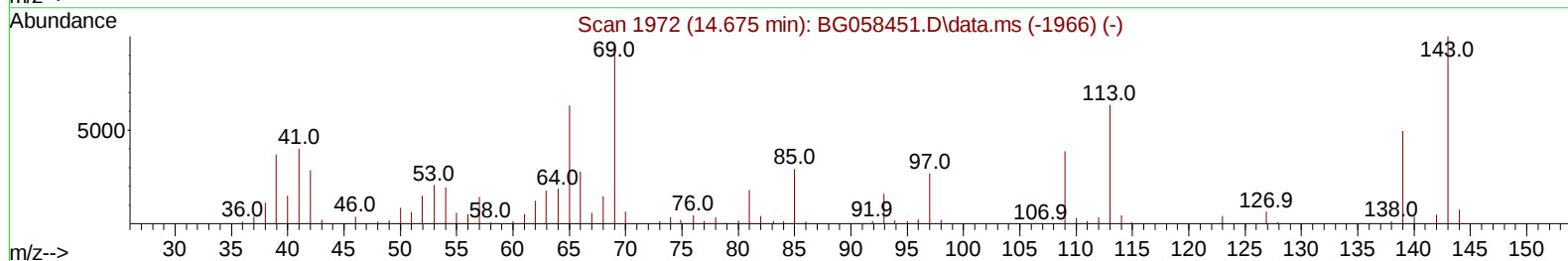
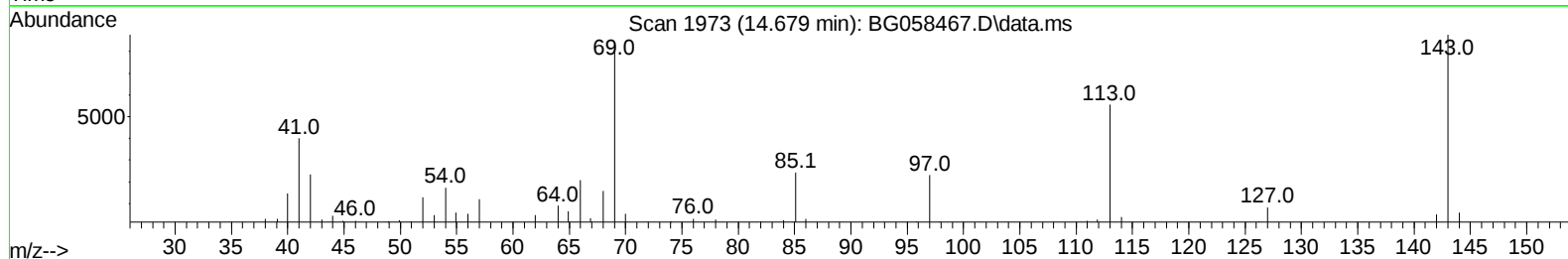
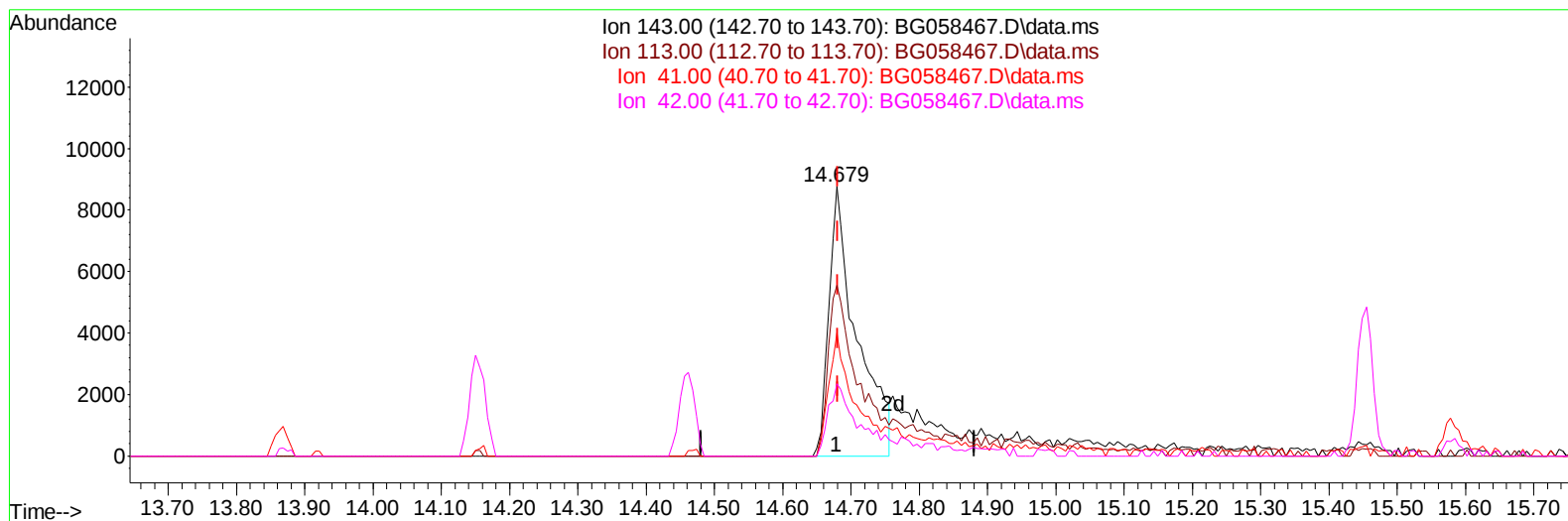
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072523\
Data File : BG058467.D
Acq On : 26 Jul 2023 6:41
Operator : CG/JU
Sample : PB154142BL
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK142

Manual IntegrationsAPPROVED

Quant Time: Jul 26 08:19:59 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 24 22:10:54 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/27/2023
Supervised By :mohammad ahmed 07/27/2023



TIC: BG058467.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.679min (-0.001) 12.48 ng/u1

response 24725

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	67.80	63.34
41.00	46.40	45.55
42.00	30.60	26.79

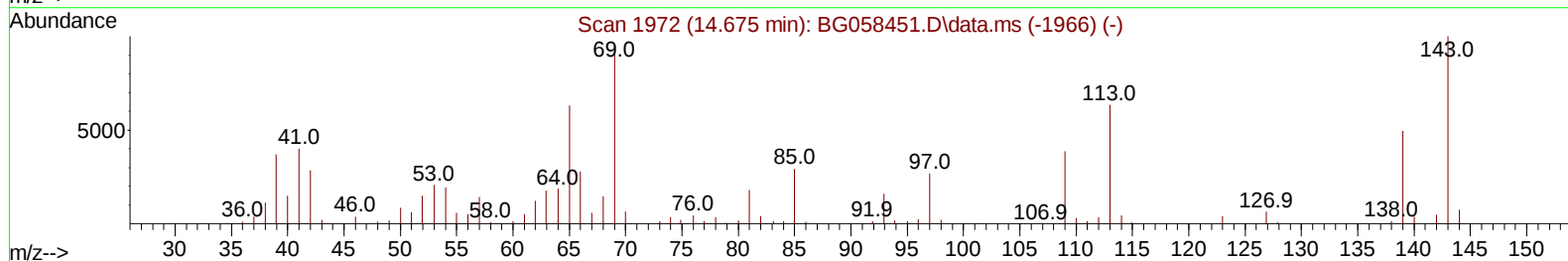
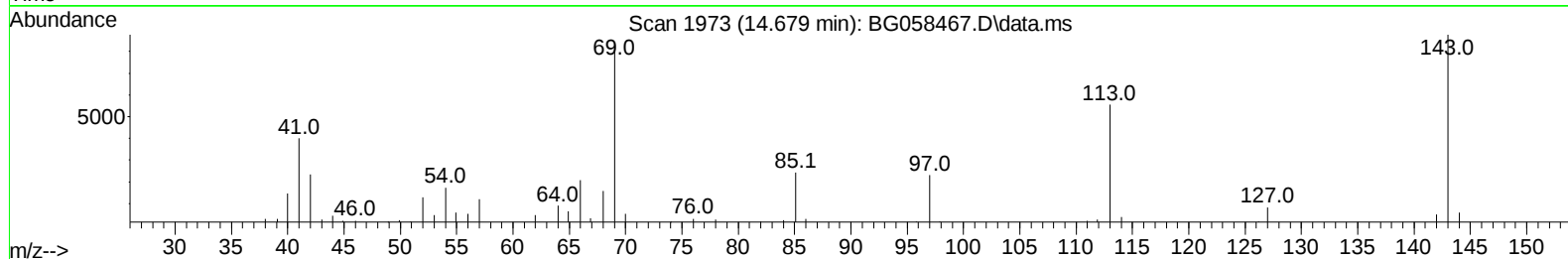
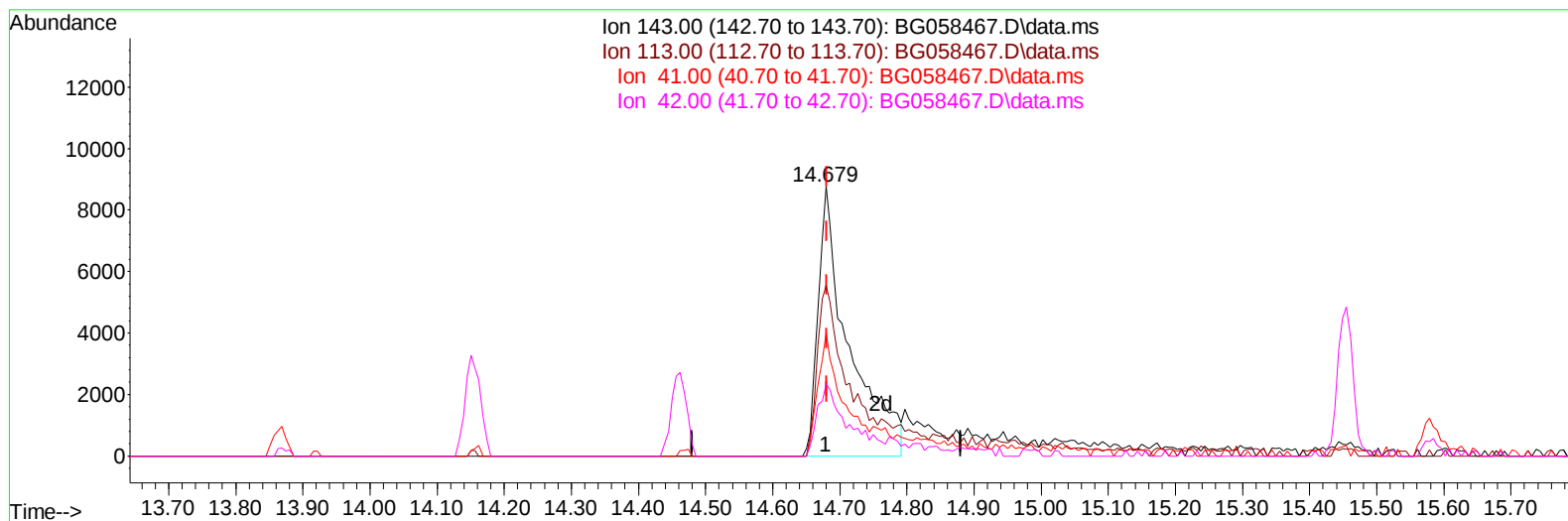
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TIC: BG058467.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.679min (-0.001) 14.05 ng/ul m

response 27833

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	67.80	63.34
41.00	46.40	45.55
42.00	30.60	26.79

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Manual IntegrationsAPPROVED

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 QLast Update : Mon Jul 24 22:10:54 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.811	152	52845	20.000	ng/ul	0.00
20) Naphthalene-d8	10.614	136	237685	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.462	164	170060	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.206	188	384566	20.000	ng/ul	0.00
79) Chrysene-d12	21.448	240	332134	20.000	ng/ul	0.00
88) Perylene-d12	24.515	264	413235	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	9928	6.915	ng/uL	0.00
4) Pyridine-d5	3.645	84	138321	32.458	ng/ul	0.00
7) Phenol-d5	6.983	99	158793	31.676	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.141	67	99916	31.889	ng/ul	0.00
11) 2-Chlorophenol-d4	7.347	132	119151	33.955	ng/ul	0.00
15) 4-Methylphenol-d8	8.528	113	116051	30.350	ng/ul	0.00
21) Nitrobenzene-d5	8.974	128	58424	32.793	ng/ul	0.00
24) 2-Nitrophenol-d4	9.703	143	66182	34.027	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.238	165	115770	31.050	ng/ul	0.00
31) 4-Chloroaniline-d4	10.755	131	155713	28.808	ng/ul	0.00
46) Dimethylphthalate-d6	13.869	166	395791	30.181	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	451066	32.822	ng/ul	0.00
54) 4-Nitrophenol-d4	14.679	143	27833m	14.054	ng/ul	0.00
60) Fluorene-d10	15.455	176	345598	31.832	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.578	200	23590	12.040	ng/ul	0.00
73) Anthracene-d10	17.306	188	544844	32.789	ng/ul	0.00
81) Pyrene-d10	19.597	212	660253	33.657	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.303	264	688882	34.491	ng/ul	0.00

Target Compounds Qvalue

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

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