

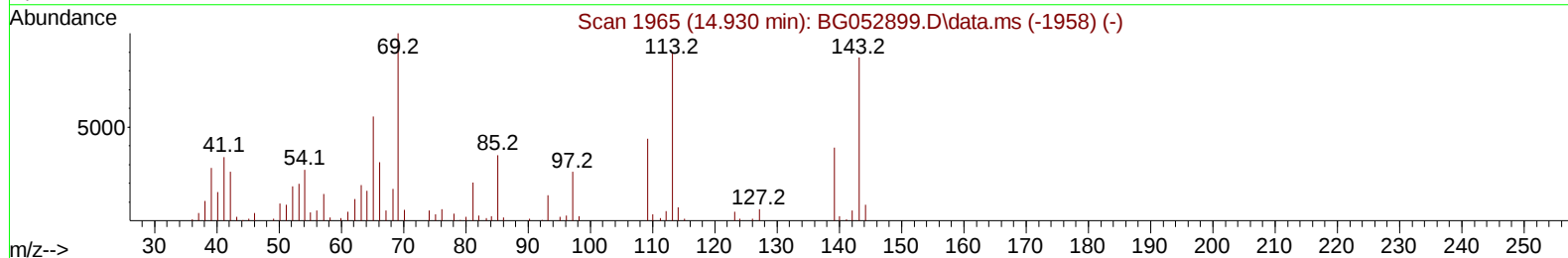
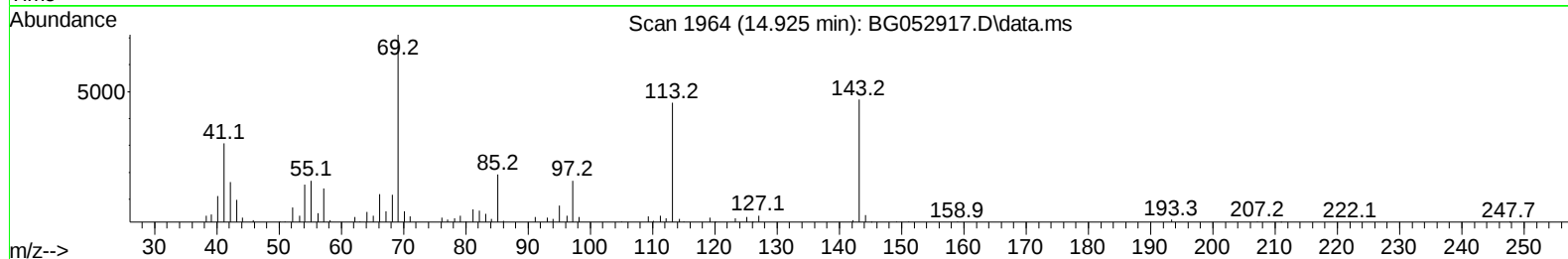
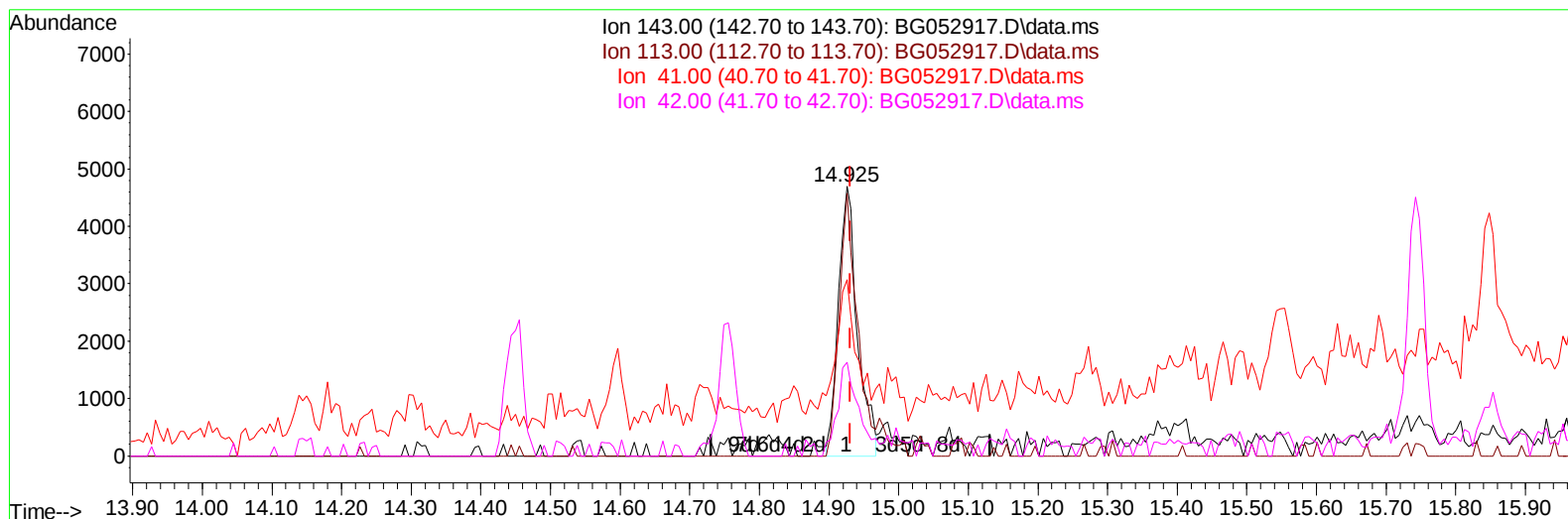
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032822\
Data File : BG052917.D
Acq On : 29 Mar 2022 8:44
Operator : CG/JU
Sample : N2097-02
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AD1

Manual IntegrationsAPPROVED

Quant Time: Mar 29 10:20:22 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Mar 23 13:16:32 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/29/2022
Supervised By :mohammad ahmed 03/29/2022



TIC: BG052917.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.925min (-0.006) 6.80 ng/u1

response 8930

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	97.51
41.00	47.20	65.45#
42.00	29.40	34.80

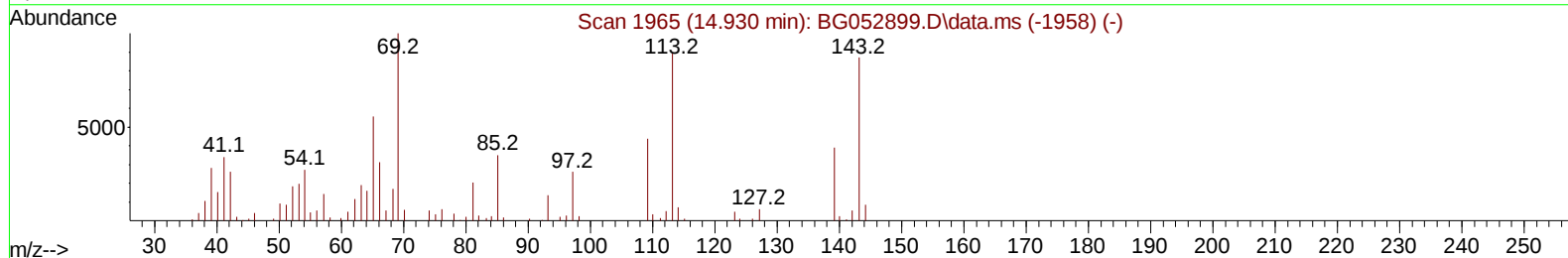
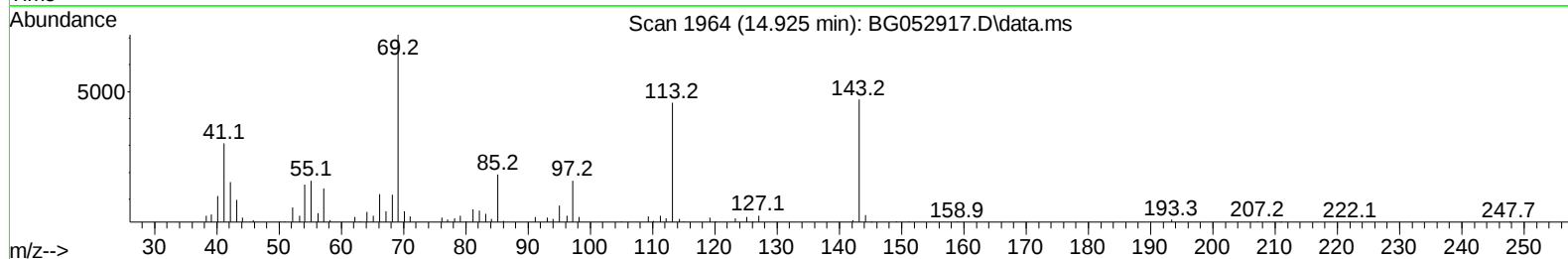
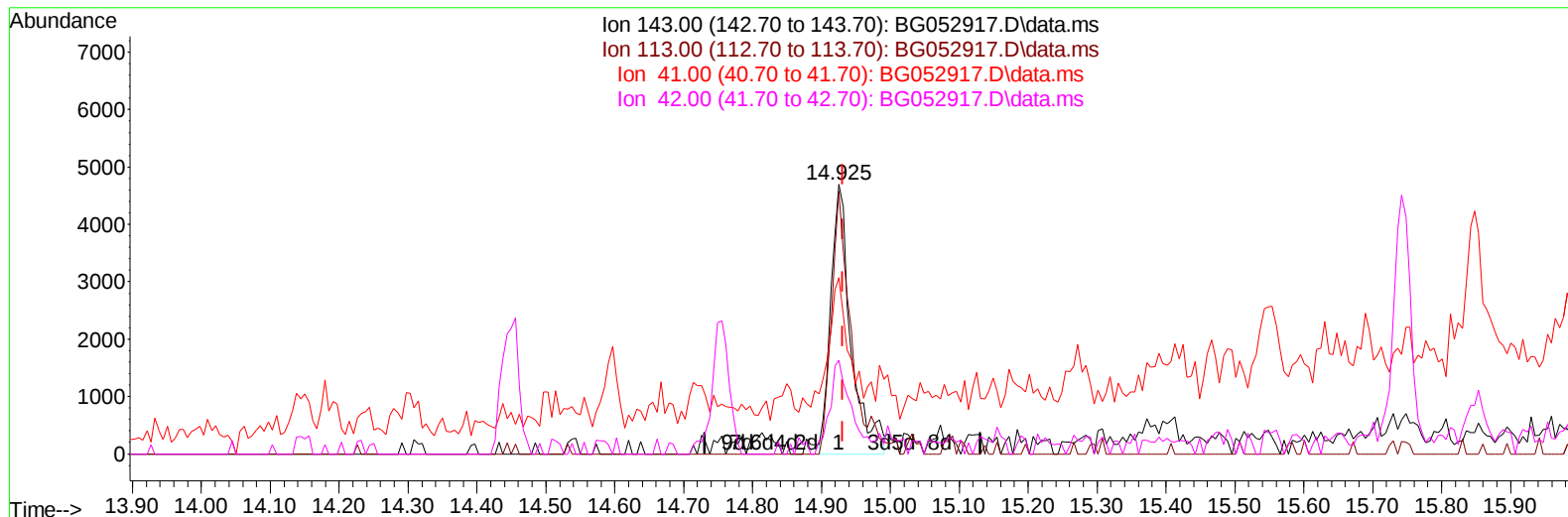
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TIC: BG052917.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.925min (-0.006) 7.30 ng/u1 m

response 9577

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	97.51
41.00	47.20	65.45#
42.00	29.40	34.80

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.135	152	27979	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.948	136	127648	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.755	164	101629	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.498	188	220076	20.000	ng/ul	0.00
79) Chrysene-d12	21.781	240	265147	20.000	ng/ul	-0.01
88) Perylene-d12	25.088	264	290386	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.547	96	3268	4.126	ng/uL	0.00
4) Pyridine-d5	3.970	84	16082	8.133	ng/ul	0.00
7) Phenol-d5	7.277	99	19729	7.873	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.453	67	54382	35.532	ng/ul	0.00
11) 2-Chlorophenol-d4	7.665	132	44078	25.904	ng/ul	-0.01
15) 4-Methylphenol-d8	8.828	113	35423	19.514	ng/ul	0.00
21) Nitrobenzene-d5	9.292	128	33519	36.479	ng/ul	0.00
24) 2-Nitrophenol-d4	10.014	143	35190	36.310	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.561	165	62969	29.865	ng/ul	0.00
31) 4-Chloroaniline-d4	11.072	131	83686	30.031	ng/ul	0.00
46) Dimethylphthalate-d6	14.150	166	276223	35.379	ng/ul	-0.01
49) Acenaphthylene-d8	14.450	160	316618	33.373	ng/ul	0.00
54) 4-Nitrophenol-d4	14.925	143	9577m	7.296	ng/ul	0.00
60) Fluorene-d10	15.742	176	242338	35.350	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.848	200	43003	38.397	ng/ul	-0.01
73) Anthracene-d10	17.598	188	382367	37.311	ng/ul	0.00
81) Pyrene-d10	19.878	212	478336	32.248	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.865	264	567382	37.815	ng/ul	-0.01

Target Compounds Qvalue

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

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