

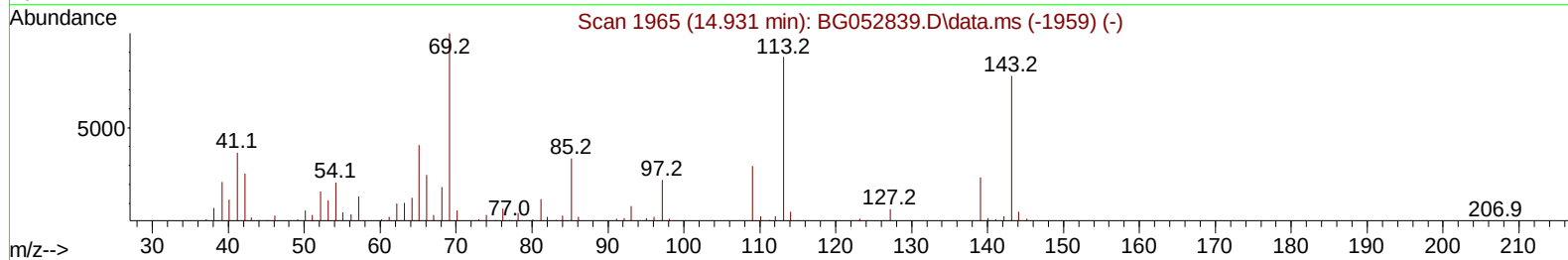
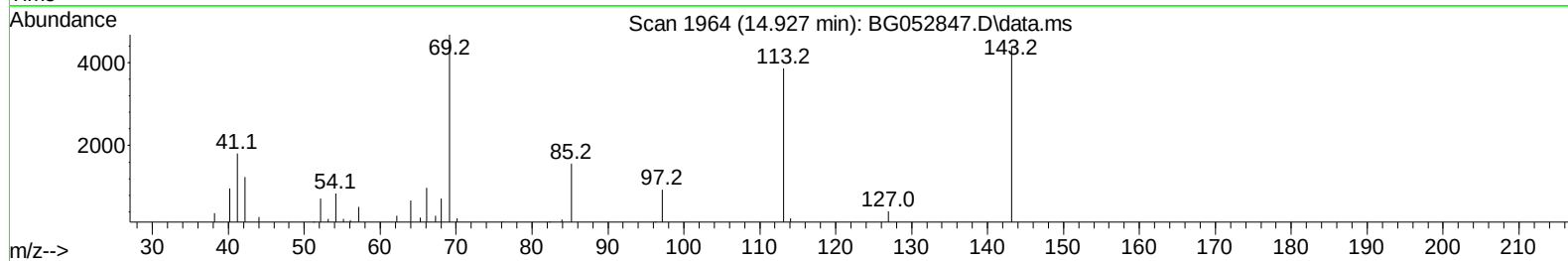
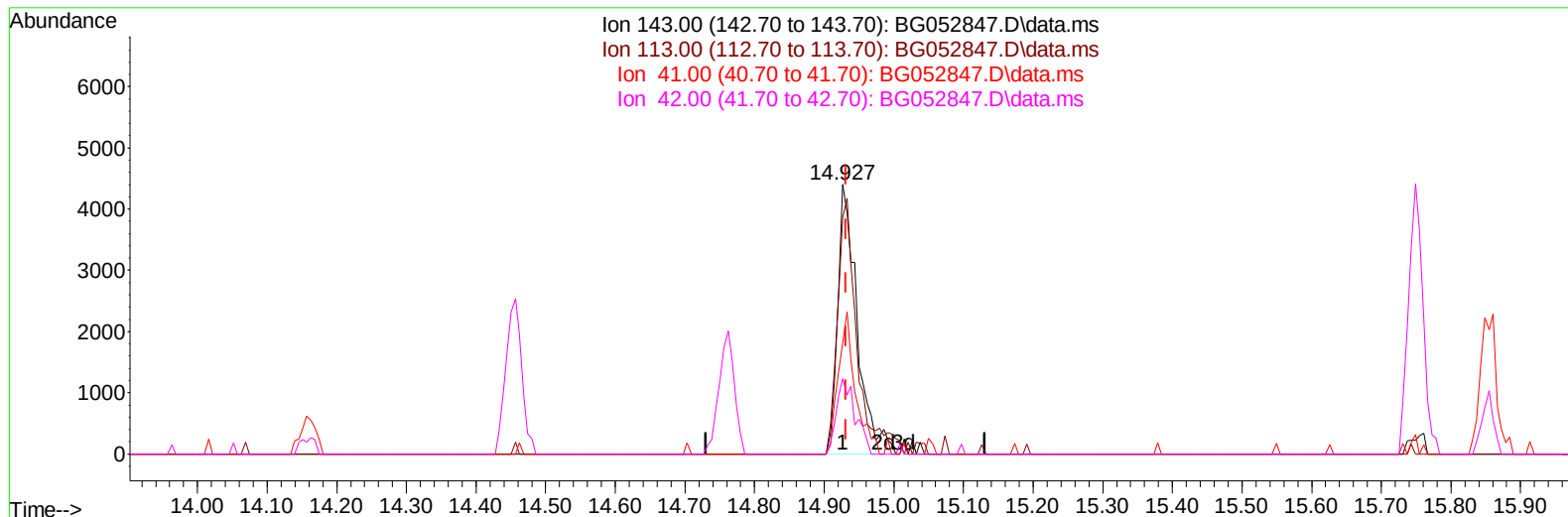
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032522\
 Data File : BG052847.D
 Acq On : 26 Mar 2022 3:06
 Operator : CG/JU
 Sample : N2079-14
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW614

Manual IntegrationsAPPROVED

Quant Time: Mar 26 05:02:44 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 03/28/2022
 Supervised By :mohammad ahmed 03/29/2022



TIC: BG052847.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.927min (-0.005) 7.10 ng/u1

response 8143

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	87.64
41.00	47.20	41.07
42.00	29.40	28.01

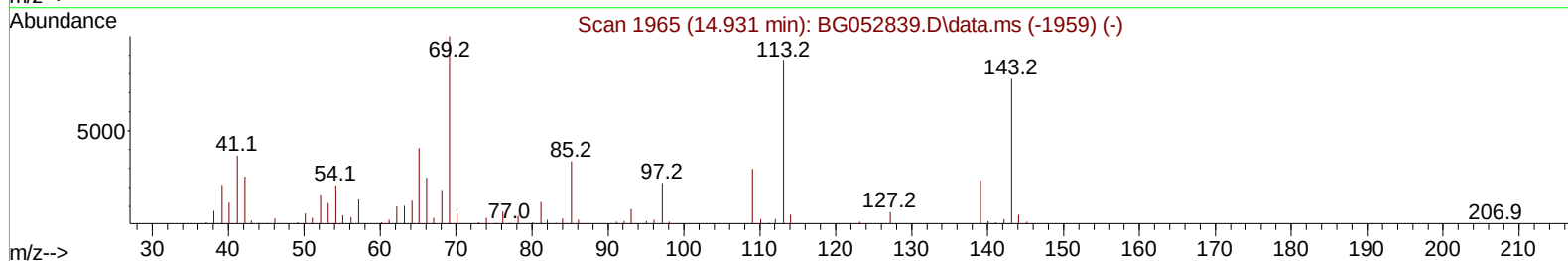
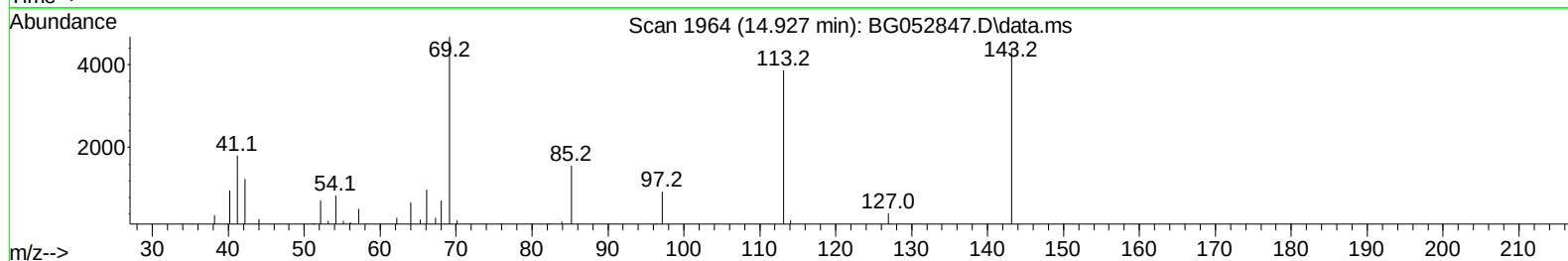
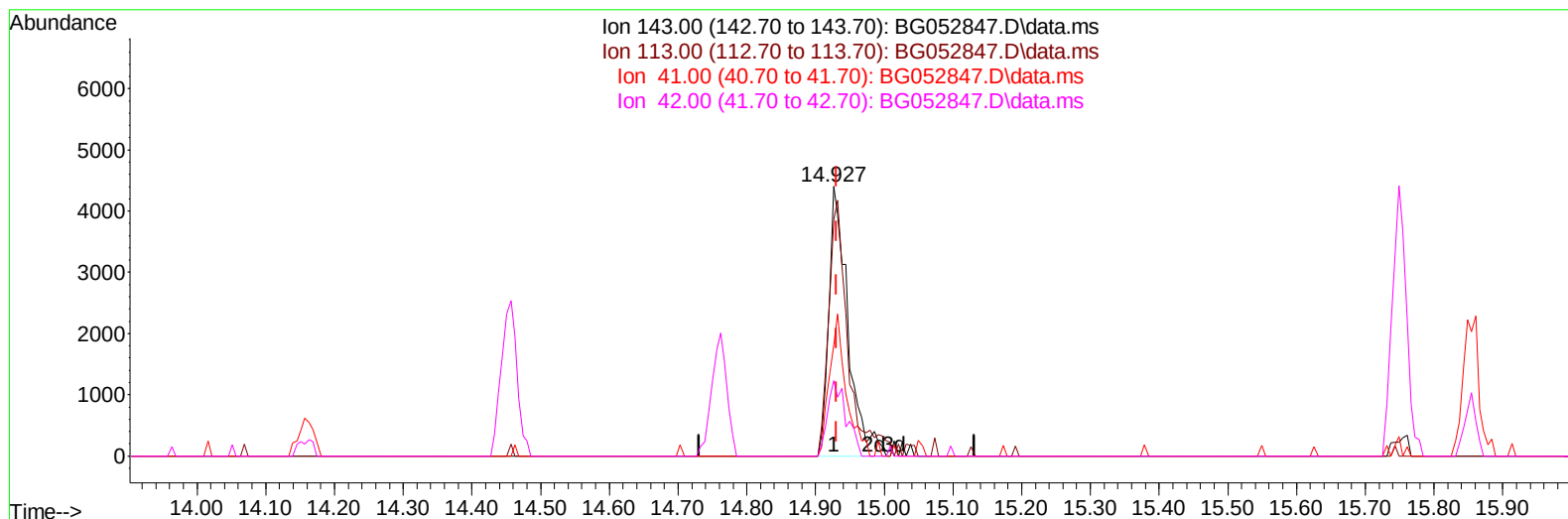
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TIC: BG052847.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.927min (-0.005) 7.43 ng/ul m

response 8530

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	87.64
41.00	47.20	41.07
42.00	29.40	28.01

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.142	152	23935	20.000	ng/ul	0.00
20) Naphthalene-d8	10.955	136	107882	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.762	164	88839	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.505	188	228222	20.000	ng/ul	0.00
79) Chrysene-d12	21.794	240	244616	20.000	ng/ul	0.00
88) Perylene-d12	25.107	264	242298	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.554	96	2848	4.203	ng/uL	0.00
4) Pyridine-d5	3.977	84	9901	5.853	ng/ul	0.00
7) Phenol-d5	7.284	99	15980	7.454	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.454	67	52040	39.746	ng/ul	0.00
11) 2-Chlorophenol-d4	7.672	132	41969	28.832	ng/ul	0.00
15) 4-Methylphenol-d8	8.835	113	28701	18.482	ng/ul	0.00
21) Nitrobenzene-d5	9.305	128	30582	39.380	ng/ul	0.00
24) 2-Nitrophenol-d4	10.027	143	34517	42.140	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.568	165	56493	31.703	ng/ul	0.00
31) 4-Chloroaniline-d4	11.079	131	69312	29.430	ng/ul	0.00
46) Dimethylphthalate-d6	14.157	166	263831	38.657	ng/ul	0.00
49) Acenaphthylene-d8	14.457	160	307064	37.025	ng/ul	0.00
54) 4-Nitrophenol-d4	14.927	143	8530m	7.434	ng/ul	0.00
60) Fluorene-d10	15.749	176	233780	39.011	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.855	200	45744	39.386	ng/ul	0.00
73) Anthracene-d10	17.605	188	434719	40.905	ng/ul	0.00
81) Pyrene-d10	19.885	212	530817	38.790	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.878	264	520898	41.607	ng/ul	0.00

Target Compounds Qvalue

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

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