

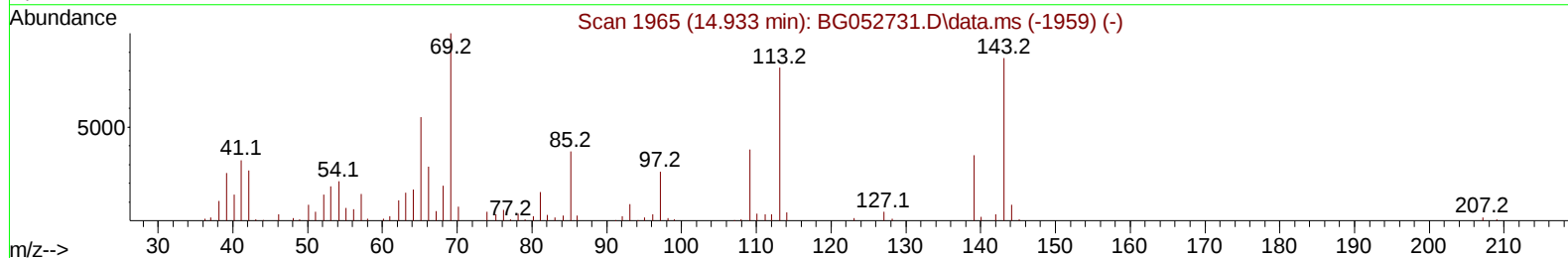
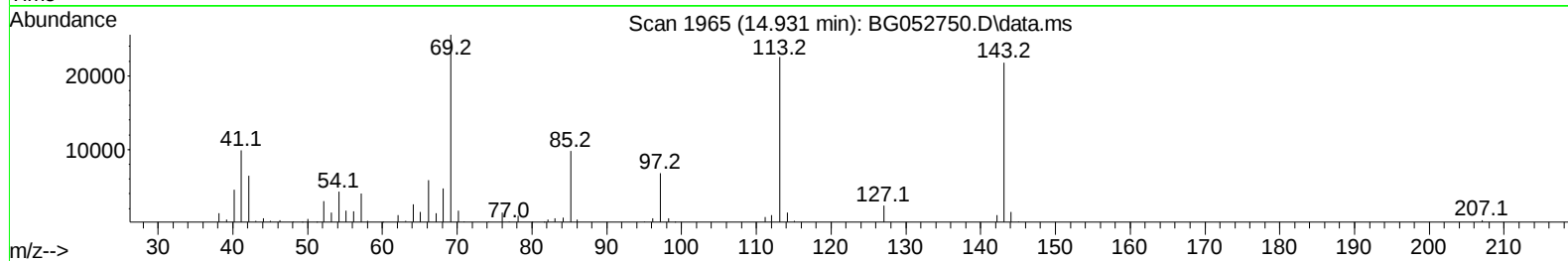
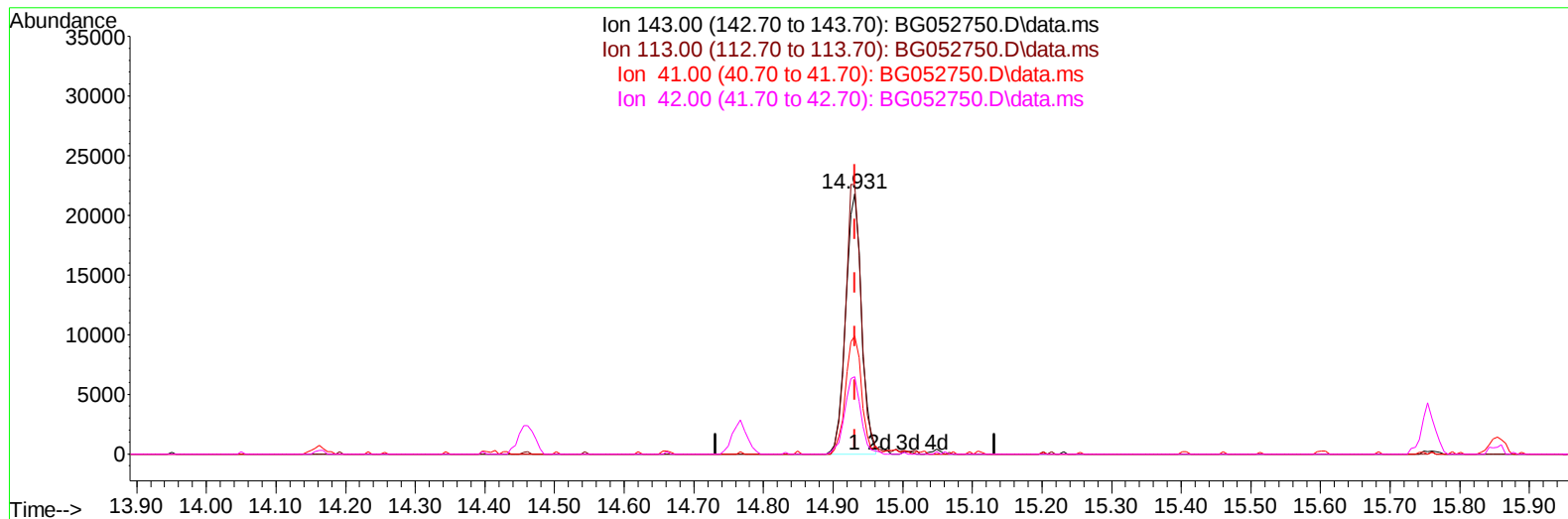
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031822\
Data File : BG052750.D
Acq On : 19 Mar 2022 18:20
Operator : CG/JU
Sample : N1820-08
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBRZ8

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 03/21/2022
Supervised By :Yogesh Patel 03/24/2022

Quant Time: Mar 21 00:51:23 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 18 00:55:42 2022
Response via : Initial Calibration



TIC: BG052750.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.931min (-0.000) 23.00 ng/u1

response 34529

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	103.37
41.00	47.20	45.23
42.00	29.40	29.66

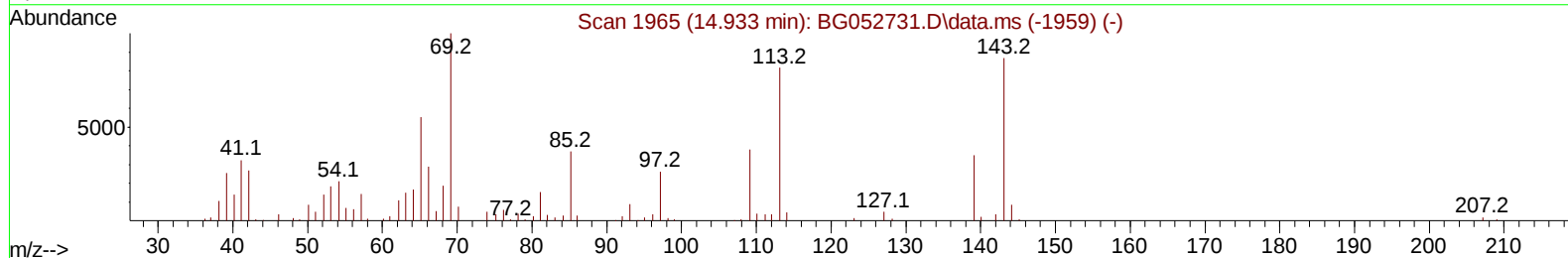
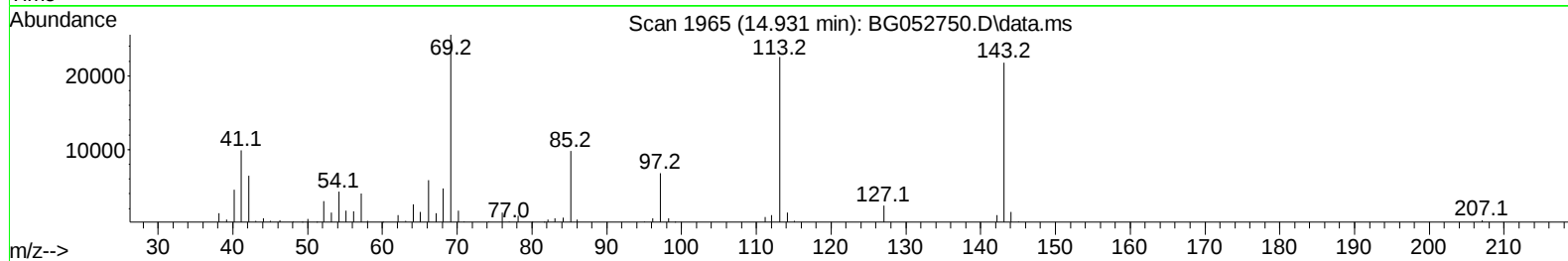
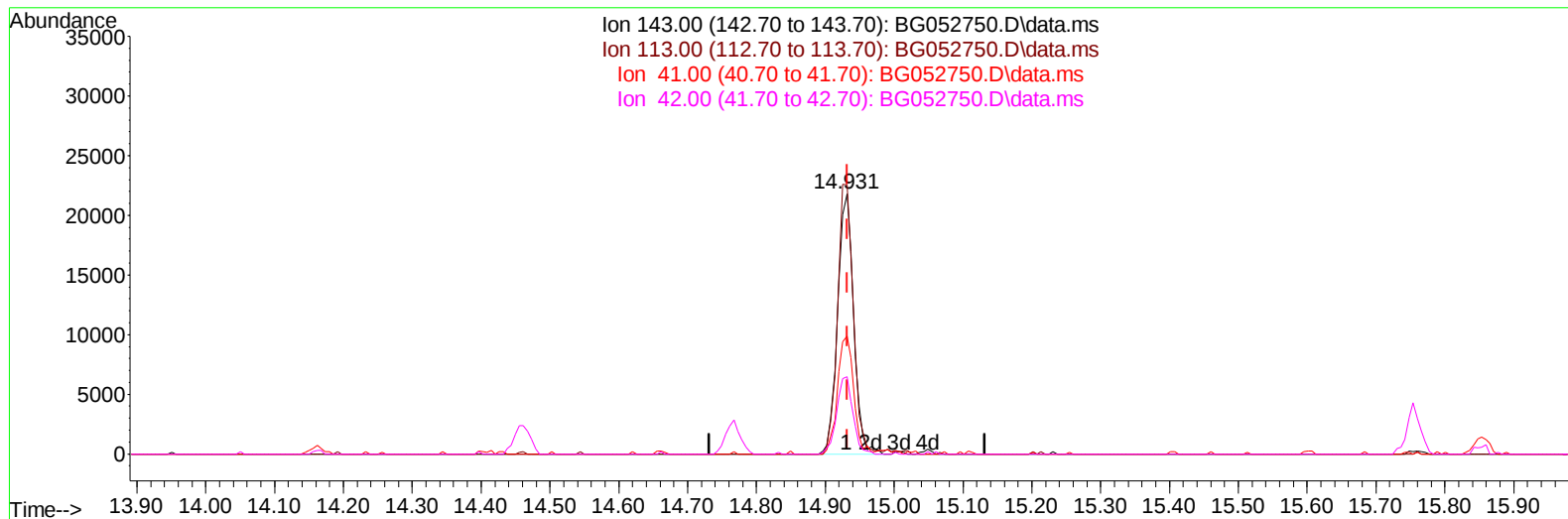
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TIC: BG052750.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.931min (-0.000) 23.31 ng/ul m

response 34991

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	103.37
41.00	47.20	45.23
42.00	29.40	29.66

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.147	152	40247	20.000	ng/ul	0.00
20) Naphthalene-d8	10.960	136	160003	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.767	164	116210	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.510	188	268637	20.000	ng/ul	0.00
79) Chrysene-d12	21.793	240	295922	20.000	ng/ul	-0.01
88) Perylene-d12	25.118	264	288072	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.559	96	5230	4.590	ng/uL	0.00
4) Pyridine-d5	3.982	84	62135	21.844	ng/ul	0.00
7) Phenol-d5	7.289	99	87813	24.360	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.465	67	56782	25.791	ng/ul	0.00
11) 2-Chlorophenol-d4	7.677	132	61974	25.320	ng/ul	0.00
15) 4-Methylphenol-d8	8.840	113	67632	25.900	ng/ul	0.00
21) Nitrobenzene-d5	9.304	128	30536	26.512	ng/ul	0.00
24) 2-Nitrophenol-d4	10.032	143	35204	28.979	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.573	165	64799	24.519	ng/ul	0.00
31) 4-Chloroaniline-d4	11.084	131	89247	25.550	ng/ul	0.00
46) Dimethylphthalate-d6	14.162	166	223207	25.001	ng/ul	0.00
49) Acenaphthylene-d8	14.461	160	276231	25.462	ng/ul	0.00
54) 4-Nitrophenol-d4	14.931	143	34991m	23.311	ng/ul	0.00
60) Fluorene-d10	15.754	176	195944	24.996	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.854	200	29254	21.399	ng/ul	0.00
73) Anthracene-d10	17.610	188	323967	25.898	ng/ul	0.00
81) Pyrene-d10	19.889	212	415022	25.070	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.889	264	386683	25.978	ng/ul	0.00

Target Compounds Qvalue

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

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