

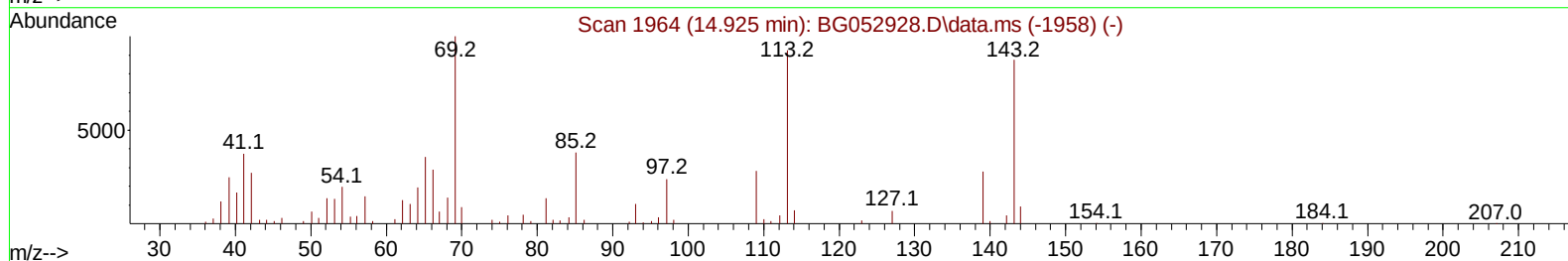
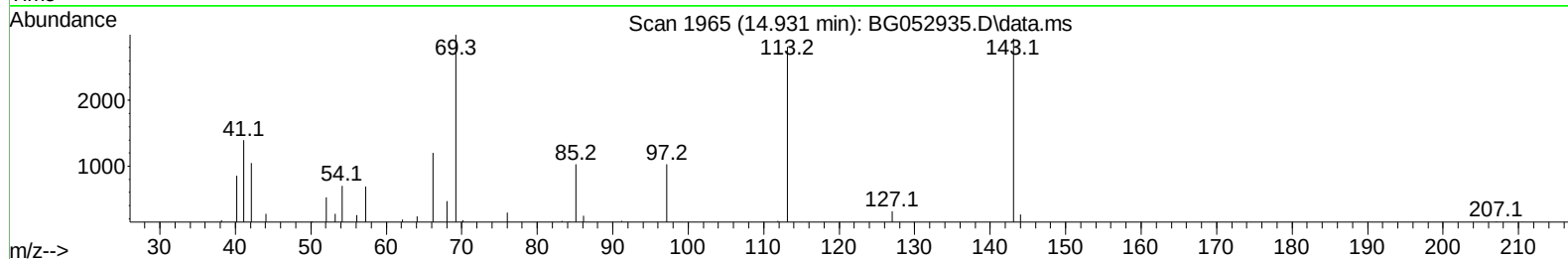
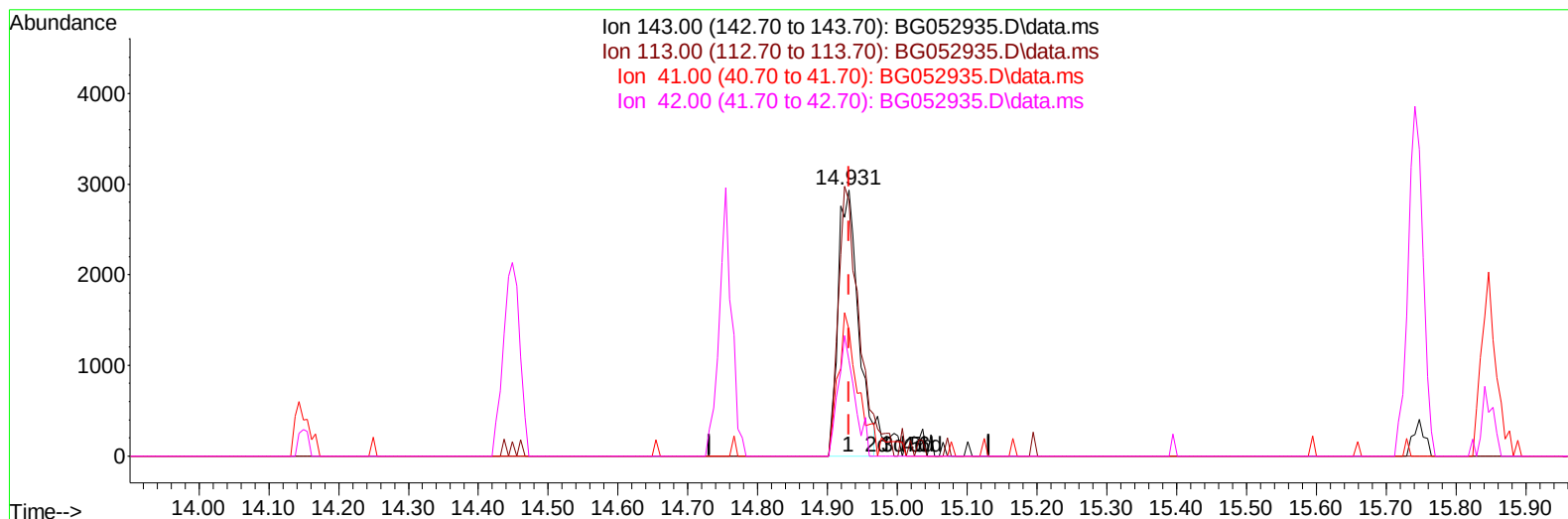
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032822\
Data File : BG052935.D
Acq On : 29 Mar 2022 20:18
Operator : CG/JU
Sample : N2097-16
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AE6

Manual IntegrationsAPPROVED

Quant Time: Mar 30 01:11:03 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 03/30/2022
Supervised By :Yogesh Patel 04/02/2022



TIC: BG052935.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.931min (-0.001) 4.79 ng/u1

response 5858

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	95.78
41.00	47.20	47.36
42.00	29.40	35.72#

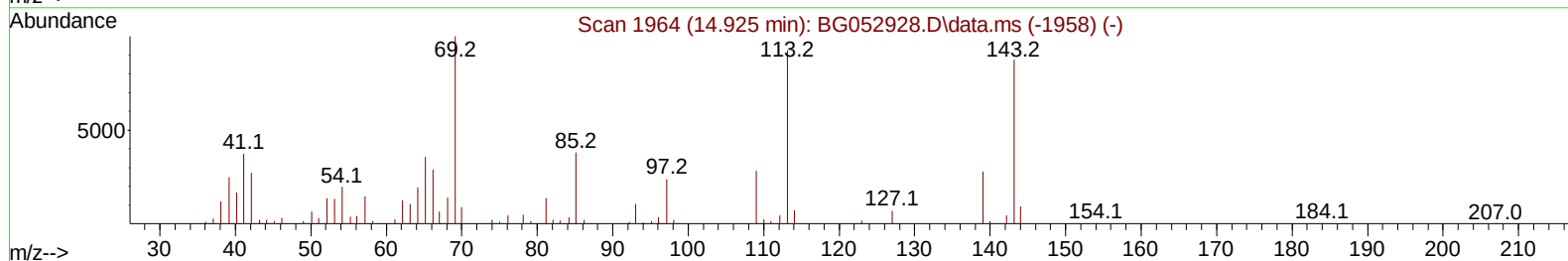
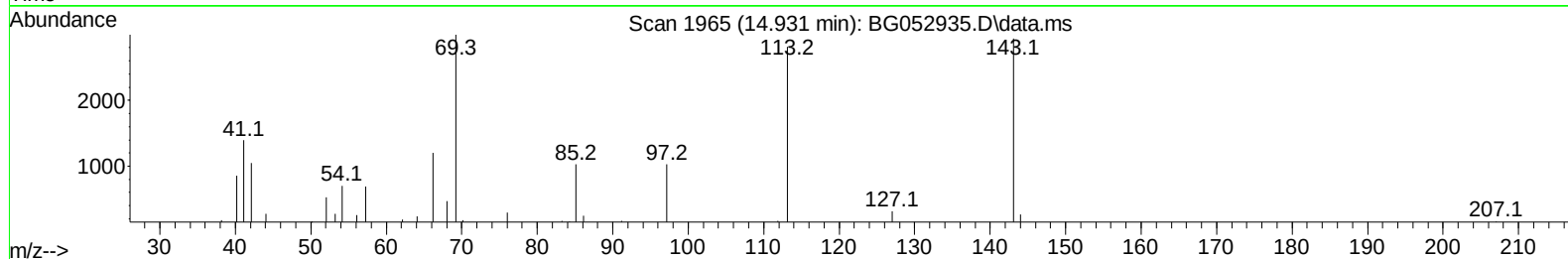
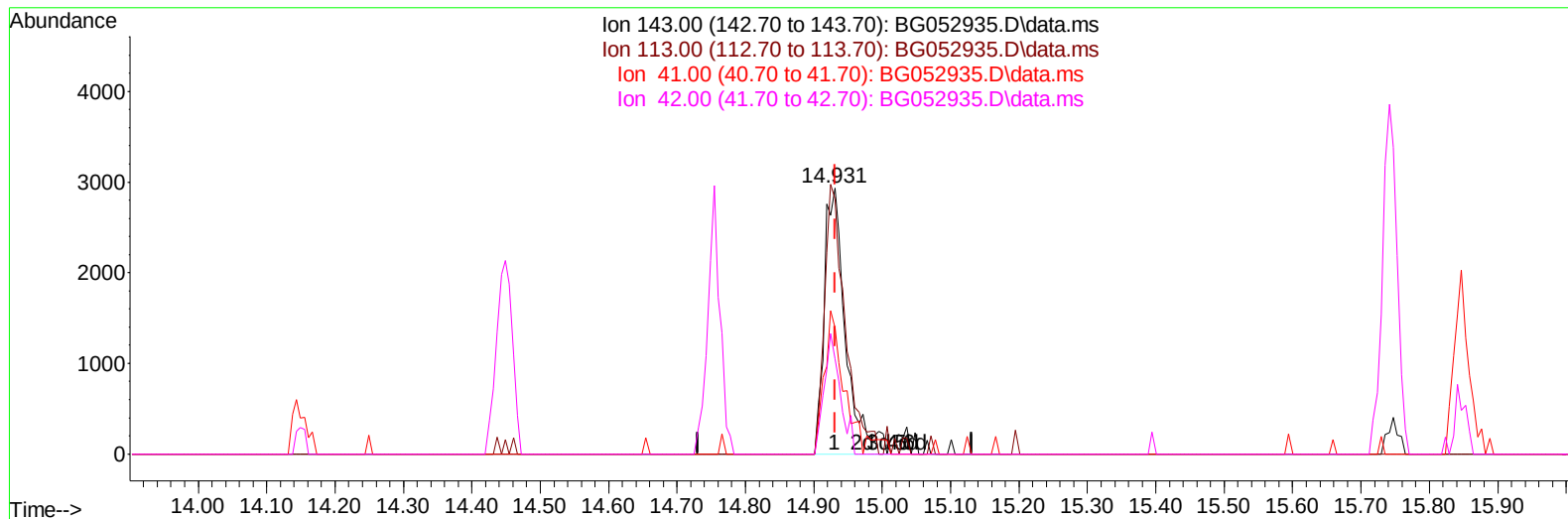
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(54) 4-Nitrophenol-d4 (S)

14.931min (-0.001) 5.22 ng/ul m

response 6388

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	95.78
41.00	47.20	47.36
42.00	29.40	35.72#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.134	152	25292	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.948	136	113855	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.754	164	94780	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.498	188	234491	20.000	ng/ul	0.00
79) Chrysene-d12	21.780	240	258459	20.000	ng/ul	#-0.01
88) Perylene-d12	25.093	264	254471	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.546	96	2630	3.673	ng/uL	0.00
4) Pyridine-d5	3.975	84	5093	2.849	ng/ul	0.00
7) Phenol-d5	7.276	99	14717	6.497	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.452	67	45338	32.770	ng/ul	0.00
11) 2-Chlorophenol-d4	7.664	132	36084	23.459	ng/ul	-0.01
15) 4-Methylphenol-d8	8.827	113	26235	15.988	ng/ul	0.00
21) Nitrobenzene-d5	9.291	128	26718	32.600	ng/ul	0.00
24) 2-Nitrophenol-d4	10.020	143	28479	32.945	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.560	165	51301	27.279	ng/ul	0.00
31) 4-Chloroaniline-d4	11.071	131	59574	23.968	ng/ul	0.00
46) Dimethylphthalate-d6	14.149	166	231980	31.859	ng/ul	-0.01
49) Acenaphthylene-d8	14.449	160	257806	29.137	ng/ul	0.00
54) 4-Nitrophenol-d4	14.931	143	6388m	5.218	ng/ul	0.00
60) Fluorene-d10	15.741	176	205633	32.163	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.847	200	36028	30.191	ng/ul	-0.01
73) Anthracene-d10	17.598	188	367497	33.656	ng/ul	0.00
81) Pyrene-d10	19.877	212	446442	30.877	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.864	264	446772	33.979	ng/ul	-0.01

Target Compounds Qvalue

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

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