

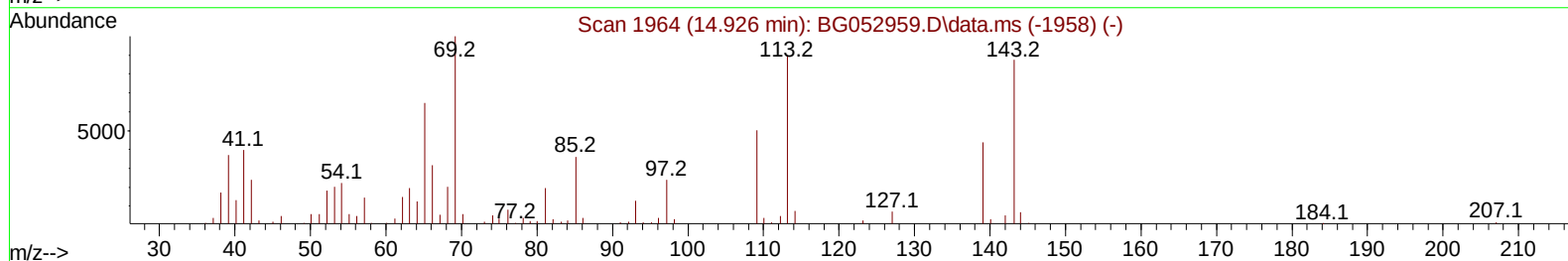
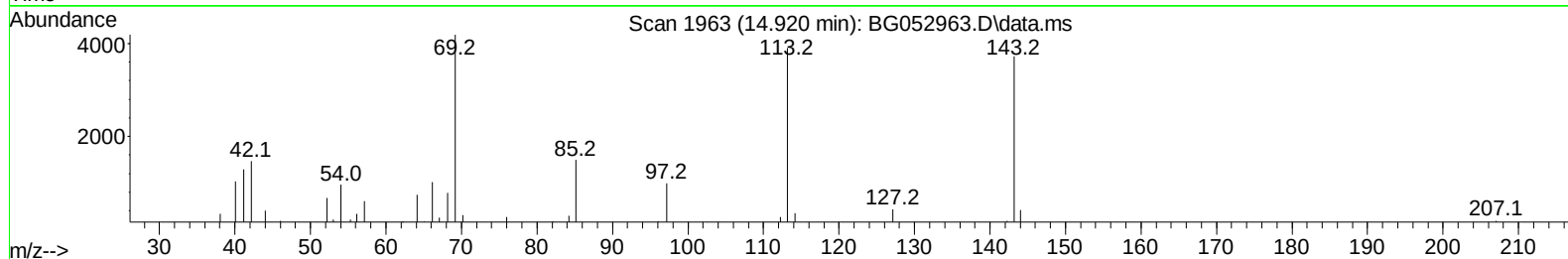
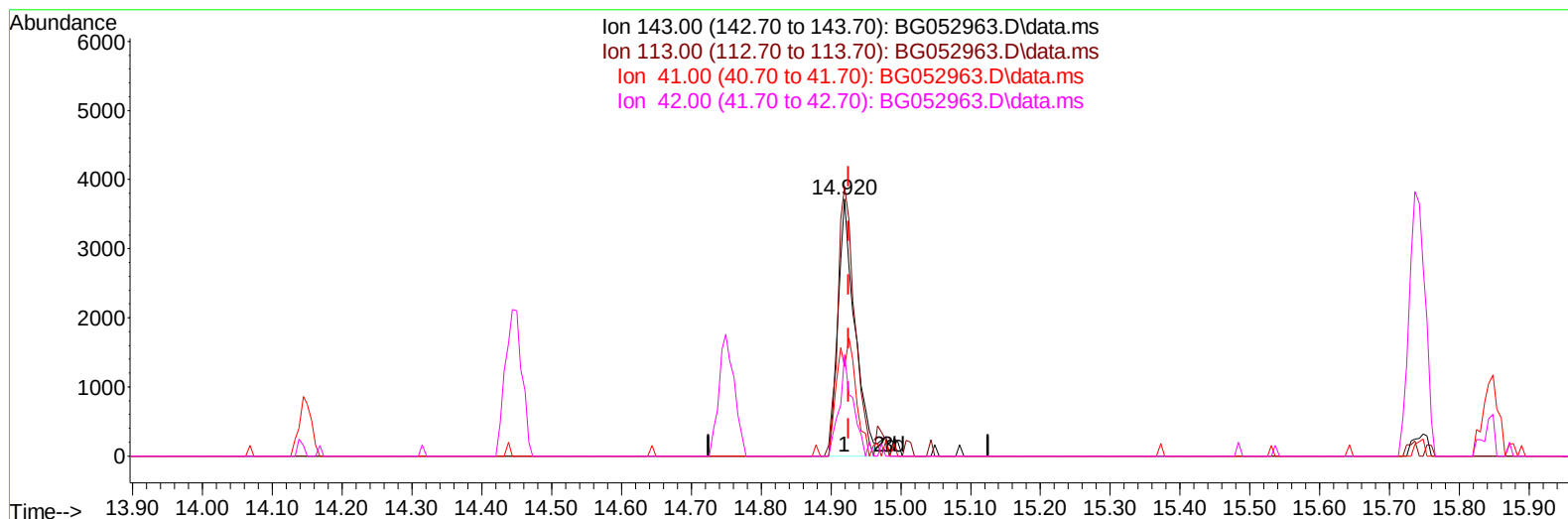
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG033022\
Data File : BG052963.D
Acq On : 31 Mar 2022 00:30
Operator : CG/JU
Sample : N2120-04
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COHT2

Manual IntegrationsAPPROVED

Quant Time: Mar 31 02:11:43 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 31 00:50:49 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/31/2022
Supervised By :Yogesh Patel 04/06/2022



TIC: BG052963.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.920min (-0.006) 5.63 ng/u1

response 6152

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	104.87
41.00	47.20	34.87#
42.00	29.40	39.39#

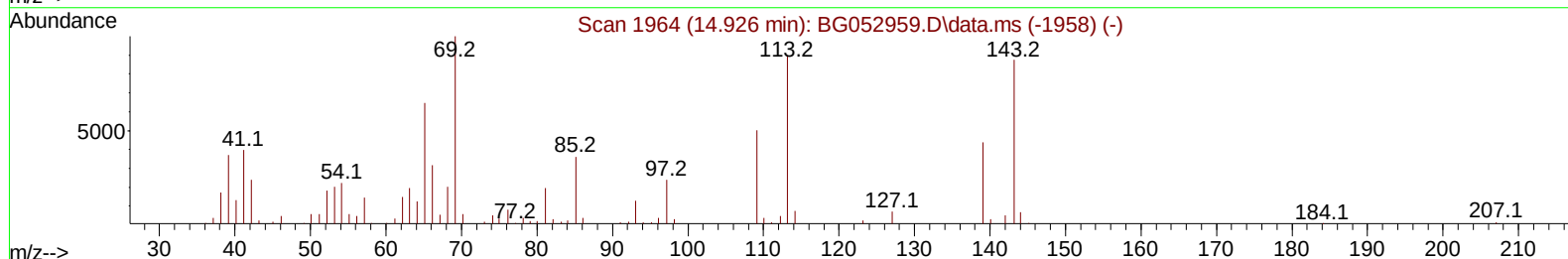
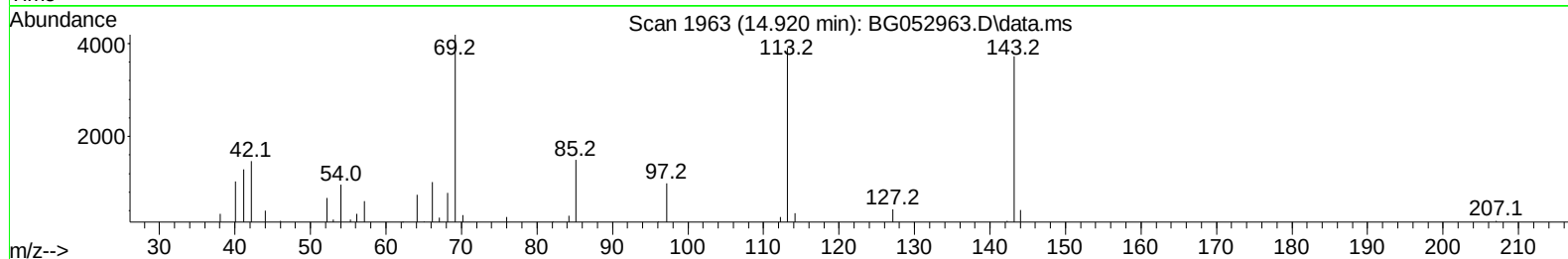
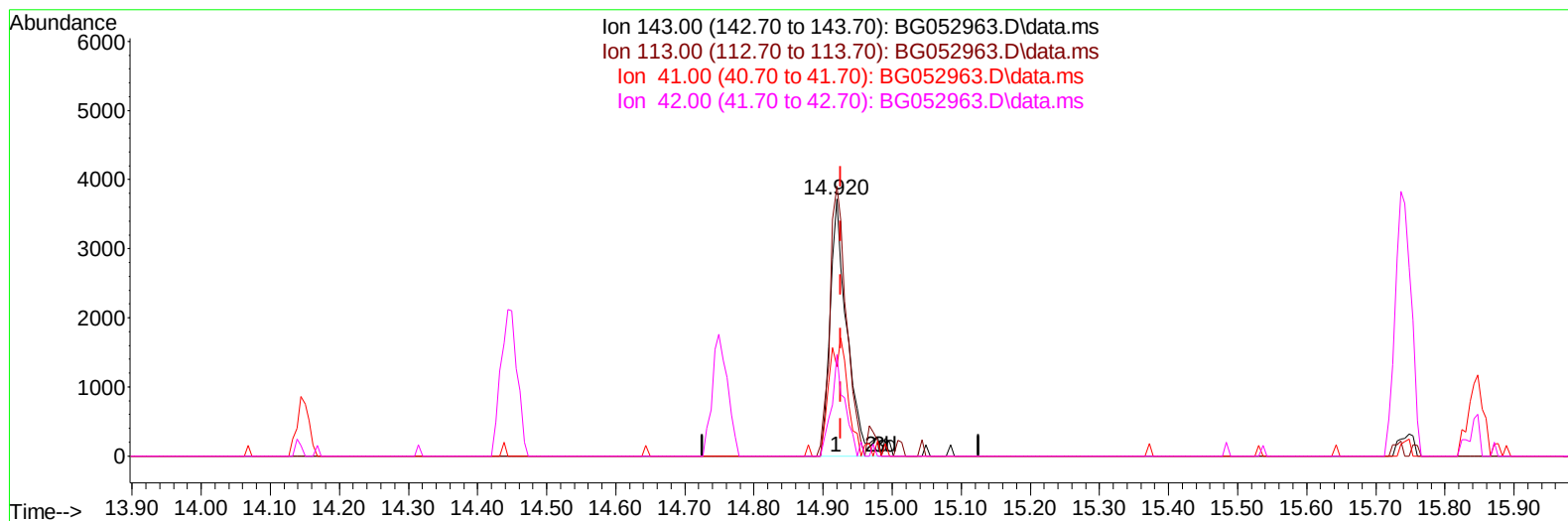
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TIC: BG052963.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.920min (-0.006) 5.85 ng/ul m

response 6395

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	104.87
41.00	47.20	34.87#
42.00	29.40	39.39#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.135	152	23973	20.000	ng/ul	0.00
20) Naphthalene-d8	10.943	136	104891	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.755	164	84643	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.493	188	209019	20.000	ng/ul	0.00
79) Chrysene-d12	21.781	240	221831	20.000	ng/ul	# 0.00
88) Perylene-d12	25.088	264	220128	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.553	96	2767	4.077	ng/uL	0.00
4) Pyridine-d5	3.976	84	9876	5.829	ng/ul	0.01
7) Phenol-d5	7.277	99	15907	7.408	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.448	67	50141	38.235	ng/ul	0.00
11) 2-Chlorophenol-d4	7.665	132	39573	27.143	ng/ul	0.00
15) 4-Methylphenol-d8	8.828	113	29565	19.008	ng/ul	0.00
21) Nitrobenzene-d5	9.292	128	29030	38.448	ng/ul	0.00
24) 2-Nitrophenol-d4	10.021	143	18806	23.614	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.555	165	50770	29.304	ng/ul	0.00
31) 4-Chloroaniline-d4	11.066	131	42277	18.463	ng/ul	0.00
46) Dimethylphthalate-d6	14.150	166	258619	39.771	ng/ul	0.00
49) Acenaphthylene-d8	14.450	160	275354	34.848	ng/ul	0.00
54) 4-Nitrophenol-d4	14.920	143	6395m	5.849	ng/ul	0.00
60) Fluorene-d10	15.742	176	230972	40.453	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.848	200	23781	22.357	ng/ul	0.00
73) Anthracene-d10	17.593	188	403216	41.427	ng/ul	0.00
81) Pyrene-d10	19.878	212	505183	40.709	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.859	264	488737	42.969	ng/ul	0.00

Target Compounds Qvalue

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

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Manual IntegrationsAPPROVED

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