

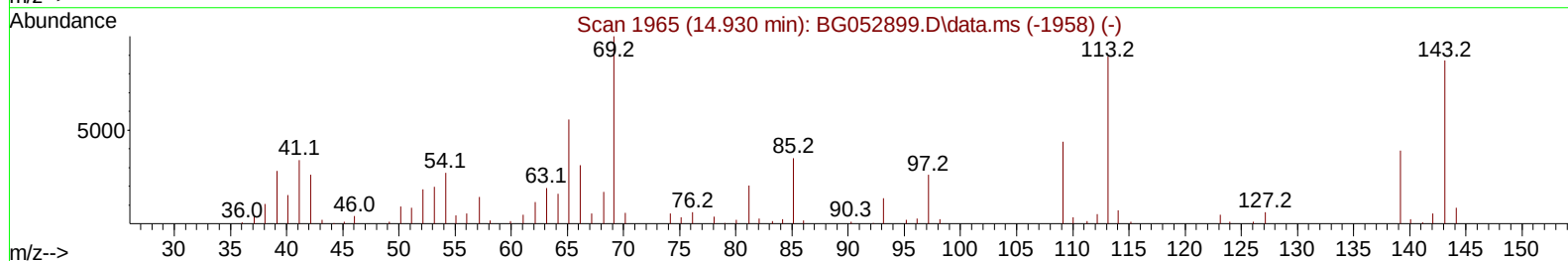
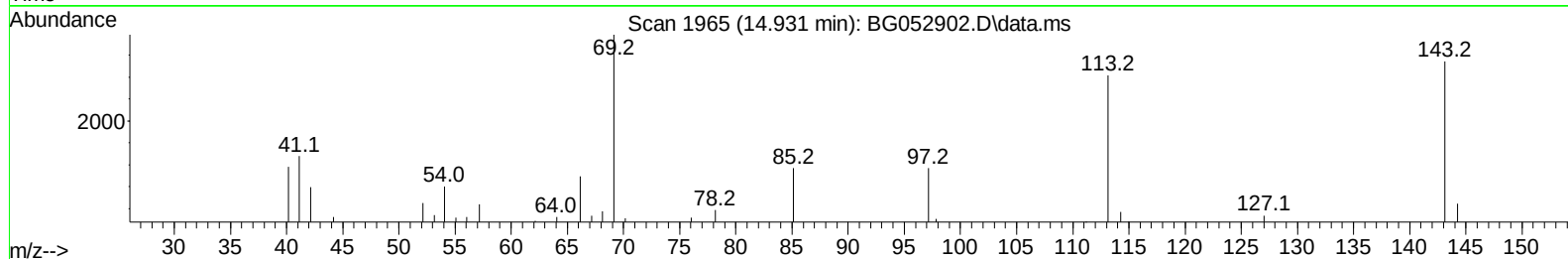
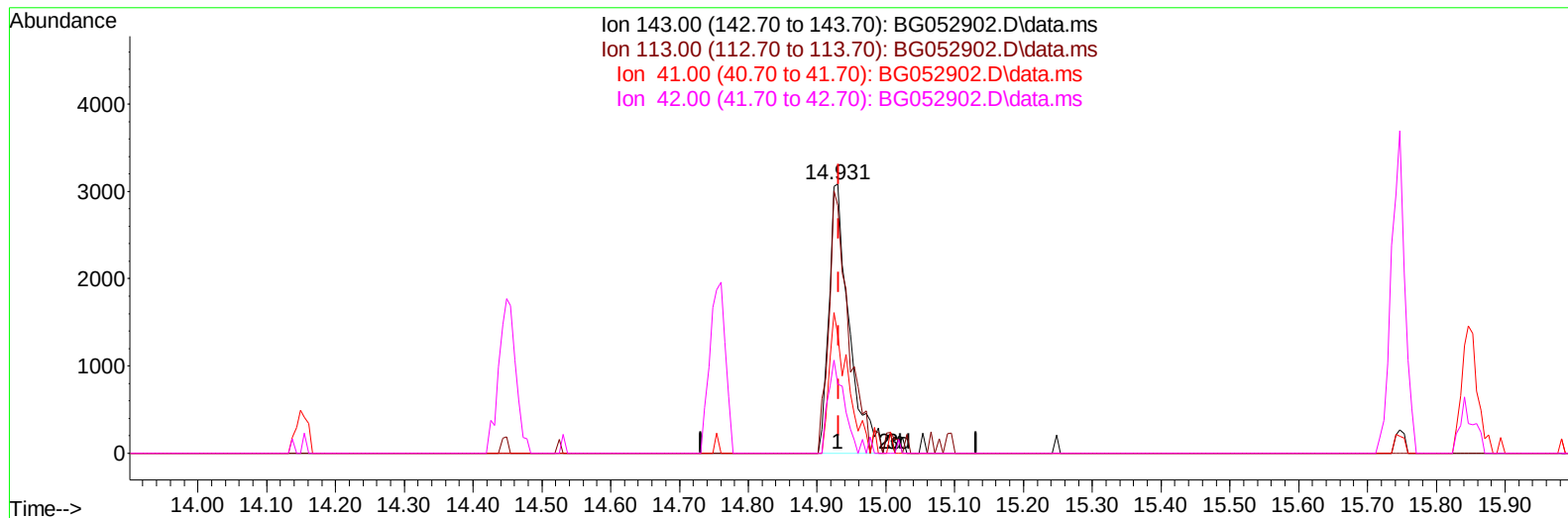
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032822\
Data File : BG052902.D
Acq On : 28 Mar 2022 22:31
Operator : CG/JU
Sample : N2100-01
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AF2

Manual IntegrationsAPPROVED

Quant Time: Mar 29 02:36:10 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: BG052902.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.931min (-0.001) 4.99 ng/ul m

response 6259

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	91.99
41.00	47.20	43.94
42.00	29.40	25.58

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.134	152	25885	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.948	136	116672	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.754	164	97085	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.498	188	233754	20.000	ng/u1	0.00
79) Chrysene-d12	21.786	240	256791	20.000	ng/u1	# 0.00
88) Perylene-d12	25.093	264	258716	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.552	96	2627	3.585	ng/uL	0.00
4) Pyridine-d5	3.969	84	16505	9.022	ng/u1	0.00
7) Phenol-d5	7.276	99	13167	5.679	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.452	67	35023	24.734	ng/u1	0.00
11) 2-Chlorophenol-d4	7.664	132	29599	18.802	ng/u1	-0.01
15) 4-Methylphenol-d8	8.833	113	22522	13.411	ng/u1	0.00
21) Nitrobenzene-d5	9.297	128	21616	25.738	ng/u1	0.00
24) 2-Nitrophenol-d4	10.020	143	23010	25.976	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.560	165	41351	21.457	ng/u1	0.00
31) 4-Chloroaniline-d4	11.071	131	67101	26.345	ng/u1	0.00
46) Dimethylphthalate-d6	14.149	166	185391	24.856	ng/u1	-0.01
49) Acenaphthylene-d8	14.449	160	206272	22.759	ng/u1	0.00
54) 4-Nitrophenol-d4	14.931	143	6259m	4.991	ng/u1	0.00
60) Fluorene-d10	15.747	176	165845	25.324	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.847	200	30120	25.320	ng/u1	-0.01
73) Anthracene-d10	17.597	188	297028	27.288	ng/u1	0.00
81) Pyrene-d10	19.877	212	373791	26.020	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.870	264	368552	27.570	ng/u1	0.00

Target Compounds Qvalue

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

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