

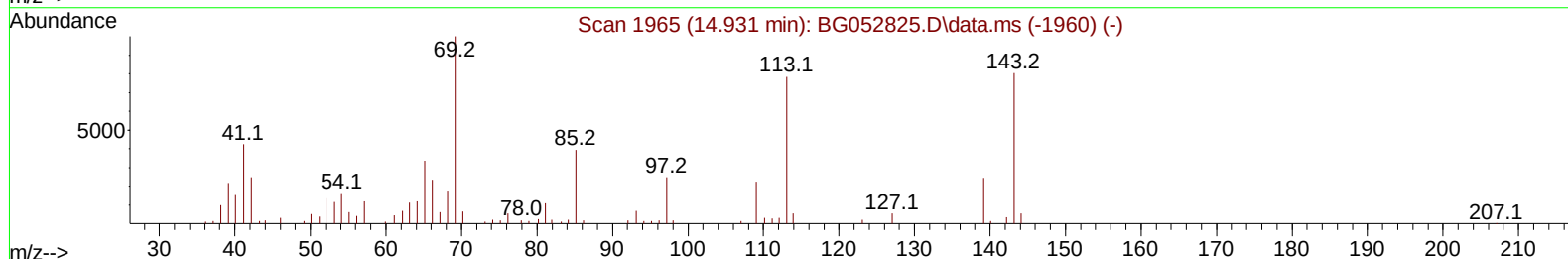
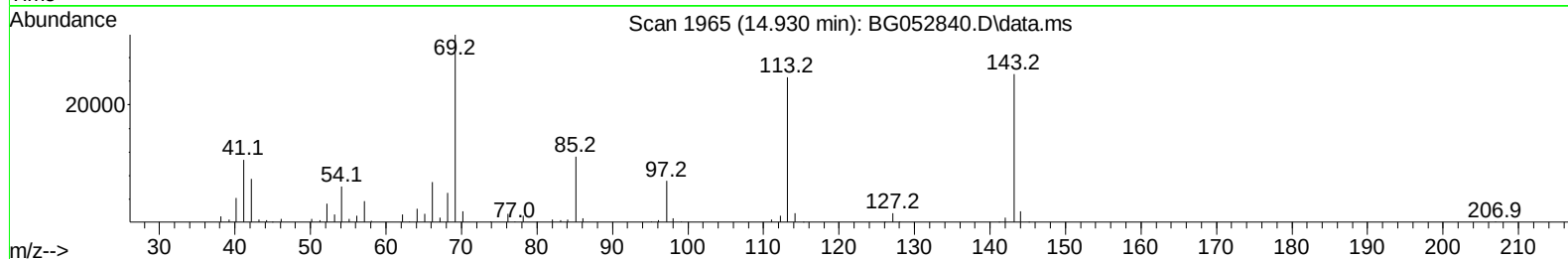
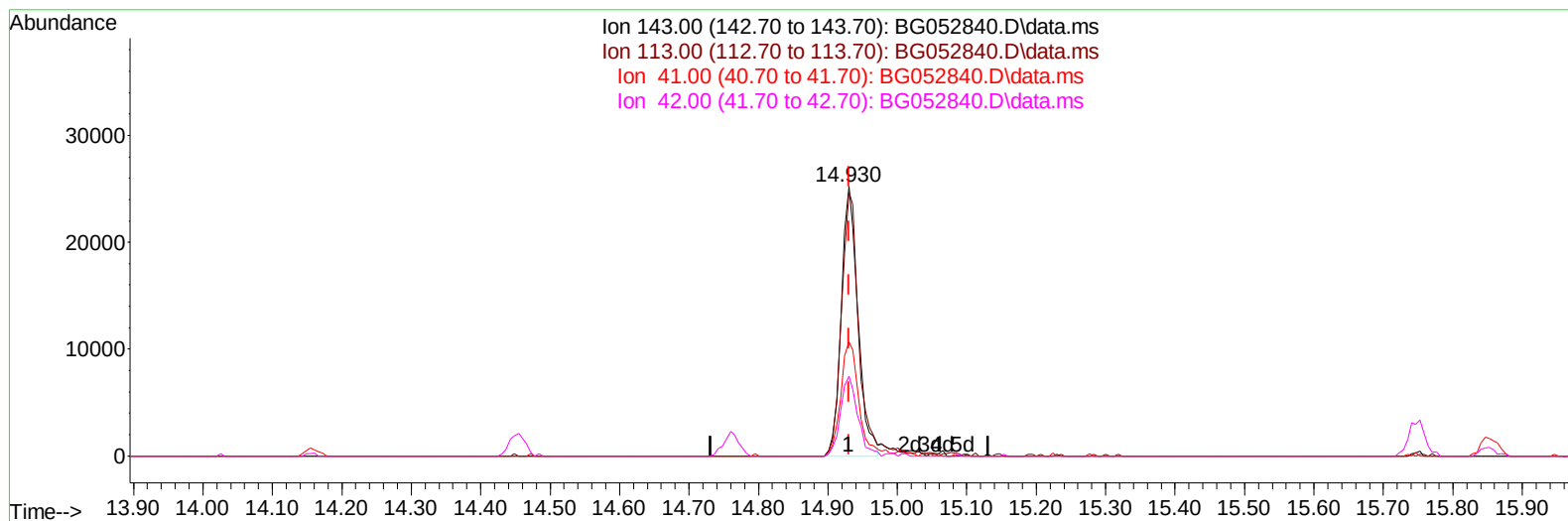
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032522\
Data File : BG052840.D
Acq On : 25 Mar 2022 22:33
Operator : CG/JU
Sample : PB143561BL
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK561

Manual IntegrationsAPPROVED

Quant Time: Mar 26 00:47:24 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: BG052840.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.930min (-0.001) 30.11 ng/ul

response 41866

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	97.88
41.00	47.20	42.39
42.00	29.40	29.71

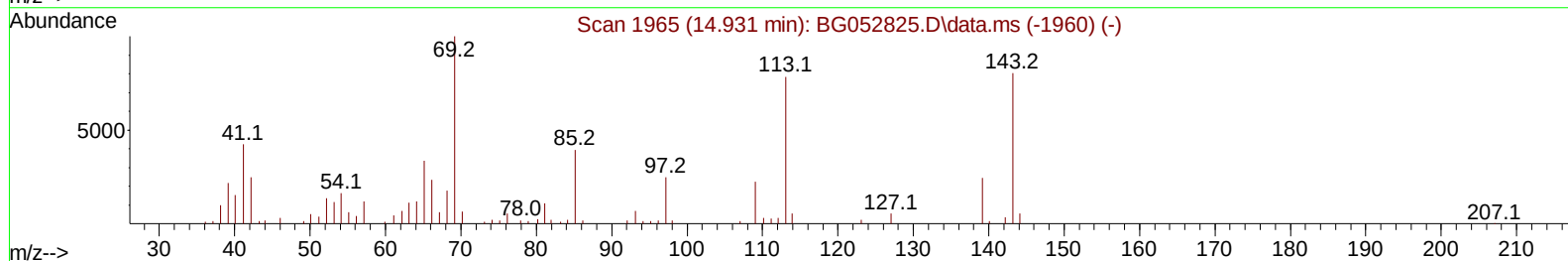
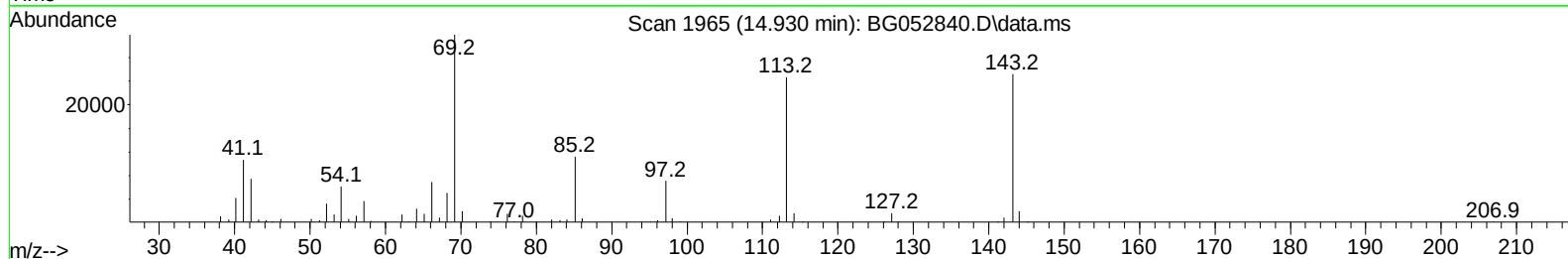
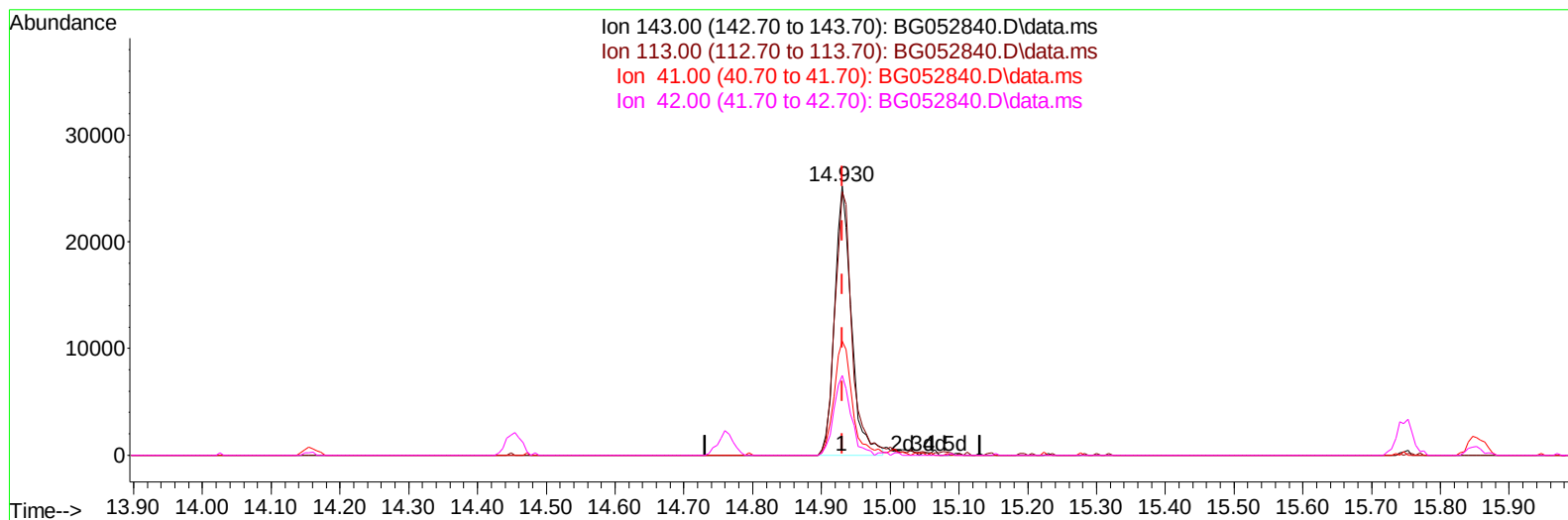
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TIC: BG052840.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.930min (-0.001) 30.77 ng/ul m

response 42793

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	97.88
41.00	47.20	42.39
42.00	29.40	29.71

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.139	152	28689	20.000	ng/ul	0.00
20) Naphthalene-d8	10.953	136	131079	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.760	164	107656	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.503	188	268456	20.000	ng/ul	0.00
79) Chrysene-d12	21.791	240	296267	20.000	ng/ul	# 0.00
88) Perylene-d12	25.110	264	283663	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.557	96	4224	5.201	ng/uL	0.00
4) Pyridine-d5	3.969	84	51448	25.373	ng/ul	0.00
7) Phenol-d5	7.282	99	75051	29.208	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.458	67	45987	29.303	ng/ul	0.00
11) 2-Chlorophenol-d4	7.669	132	49708	28.490	ng/ul	0.00
15) 4-Methylphenol-d8	8.833	113	57396	30.836	ng/ul	0.00
21) Nitrobenzene-d5	9.297	128	28487	30.191	ng/ul	0.00
24) 2-Nitrophenol-d4	10.025	143	32441	32.597	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.565	165	58332	26.942	ng/ul	0.00
31) 4-Chloroaniline-d4	11.077	131	82829	28.946	ng/ul	0.00
46) Dimethylphthalate-d6	14.155	166	251536	30.413	ng/ul	0.00
49) Acenaphthylene-d8	14.454	160	277906	27.652	ng/ul	0.00
54) 4-Nitrophenol-d4	14.930	143	42793m	30.774	ng/ul	0.00
60) Fluorene-d10	15.753	176	211231	29.087	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.852	200	39172	28.673	ng/ul	0.00
73) Anthracene-d10	17.603	188	369315	29.543	ng/ul	0.00
81) Pyrene-d10	19.882	212	460211	27.767	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.881	264	436553	29.785	ng/ul	0.00

Target Compounds Qvalue

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

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