

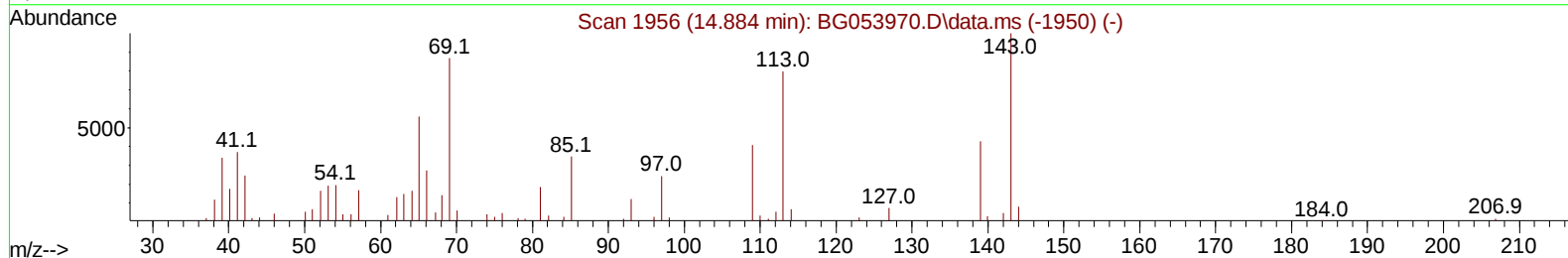
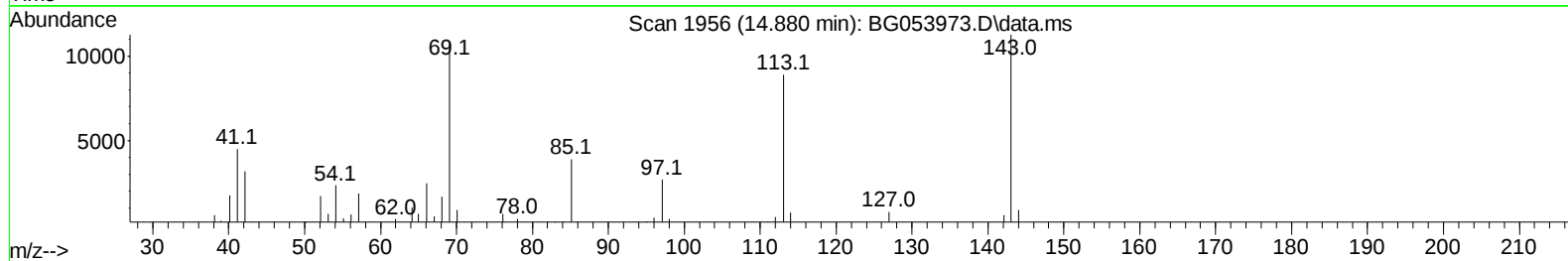
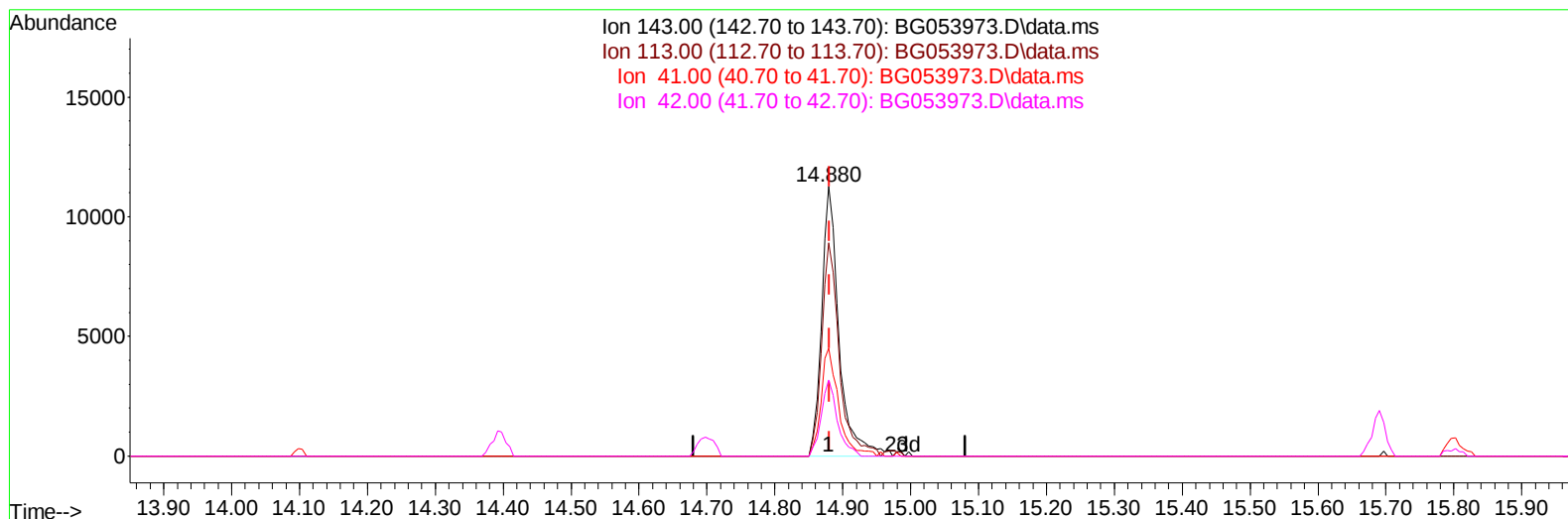
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062122\
Data File : BG053973.D
Acq On : 21 Jun 2022 11:34
Operator : CG/JU
Sample : PB145591BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK591

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 06/22/2022
Supervised By :mohammad ahmed 06/28/2022

Quant Time: Jun 21 23:06:46 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062022.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jun 20 15:29:13 2022
Response via : Initial Calibration



TIC: BG053973.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.880min (-0.001) 31.22 ng/u1****response 19910**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.90	79.13
41.00	51.40	39.95#
42.00	37.10	28.24#

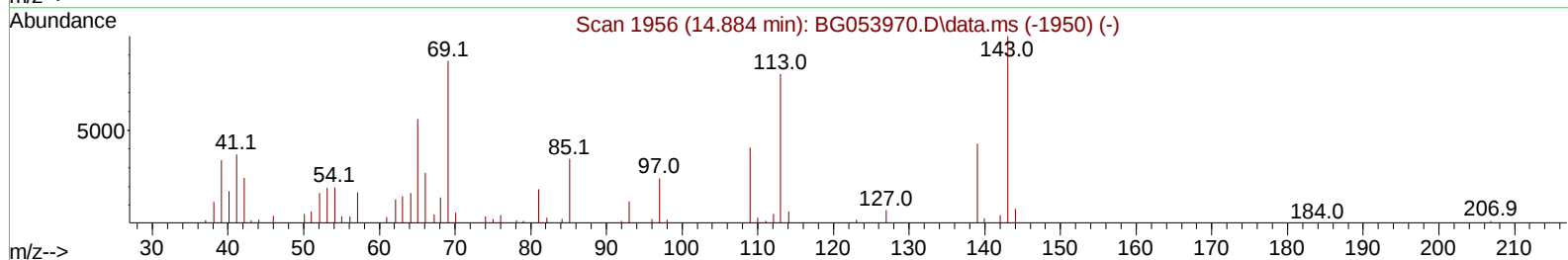
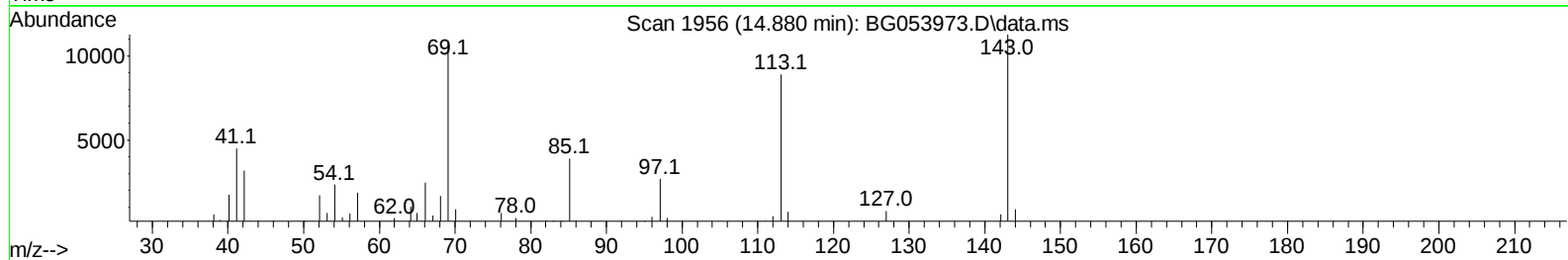
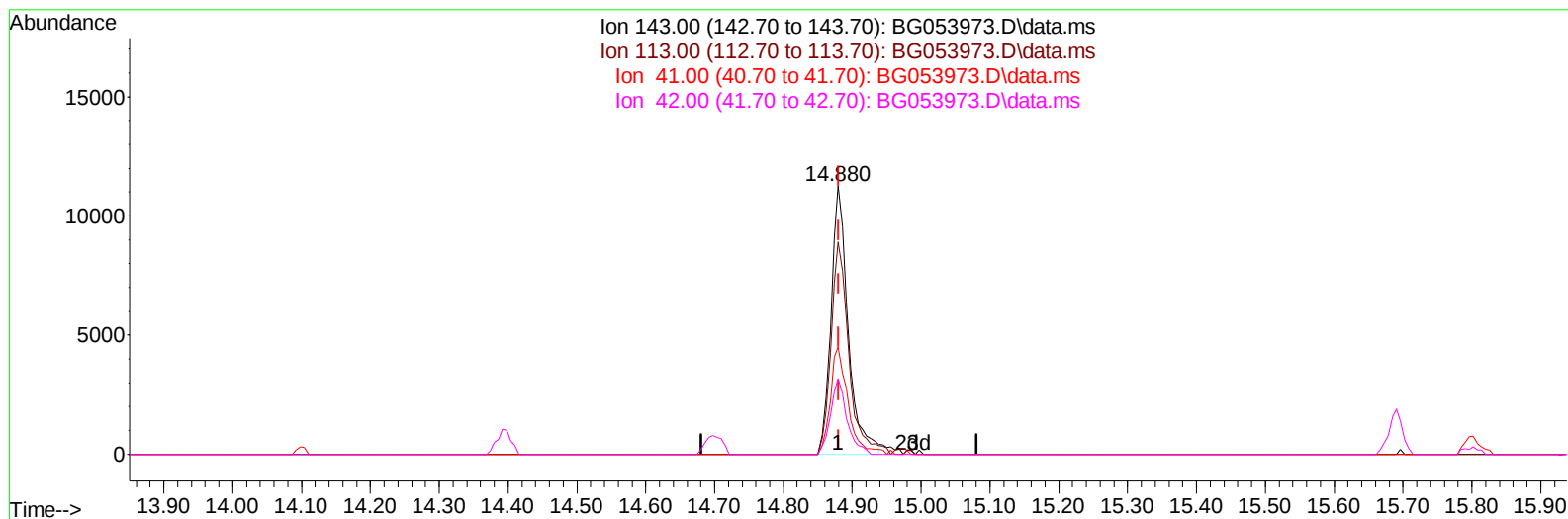
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TIC: BG053973.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.880min (-0.001) 30.43 ng/ul m

response 19410

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.90	79.13
41.00	51.40	39.95#
42.00	37.10	28.24#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.082	152	18508	20.000	ng/ul	0.00
20) Naphthalene-d8	10.884	136	74581	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.703	164	59213	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.441	188	162472	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.719	240	188997	20.000	ng/ul	# 0.00
88) Perylene-d12	24.980	264	190405	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.505	96	2430	6.247	ng/uL	0.00
4) Pyridine-d5	3.916	84	34683	32.781	ng/ul	0.00
7) Phenol-d5	7.230	99	42418	31.991	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.400	67	26612	33.272	ng/ul	0.00
11) 2-Chlorophenol-d4	7.606	132	36072	34.115	ng/ul	0.00
15) 4-Methylphenol-d8	8.775	113	36484	32.271	ng/ul	0.00
21) Nitrobenzene-d5	9.239	128	17823	32.196	ng/ul	0.00
24) 2-Nitrophenol-d4	9.962	143	20356	34.608	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.502	165	41233	31.685	ng/ul	0.00
31) 4-Chloroaniline-d4	11.014	131	52242	31.777	ng/ul	0.00
46) Dimethylphthalate-d6	14.098	166	151359	32.493	ng/ul	0.00
49) Acenaphthylene-d8	14.392	160	159639	31.516	ng/ul	0.00
54) 4-Nitrophenol-d4	14.880	143	19410m	30.431	ng/ul	0.00
60) Fluorene-d10	15.691	176	134322	32.102	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.802	200	22682	25.592	ng/ul	0.00
73) Anthracene-d10	17.541	188	242817	33.093	ng/ul	0.00
81) Pyrene-d10	19.827	212	318419	32.313	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.756	264	323581	33.667	ng/ul	0.00

Target Compounds Qvalue

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

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