

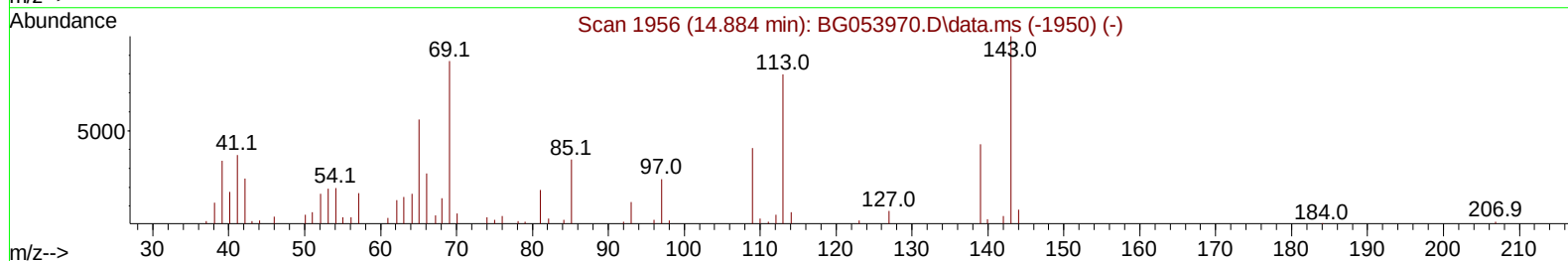
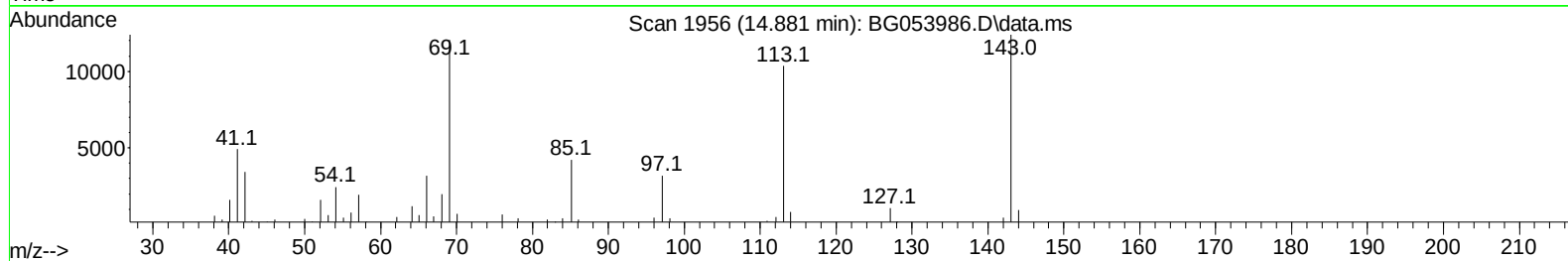
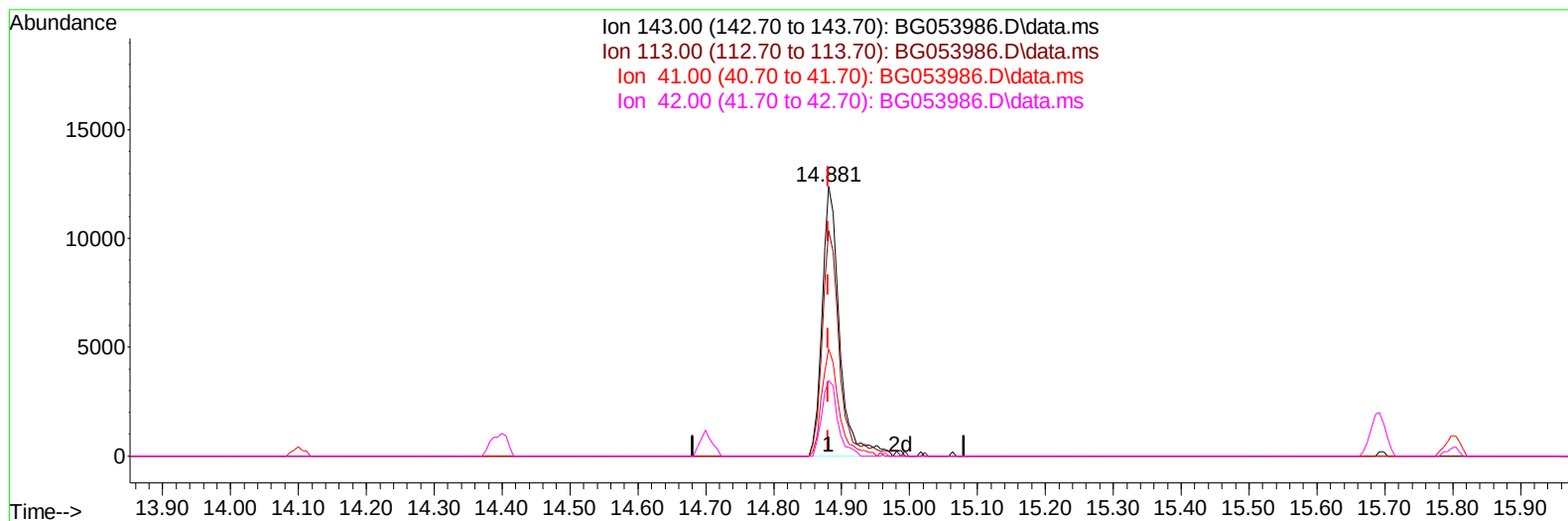
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062122\
Data File : BG053986.D
Acq On : 21 Jun 2022 20:39
Operator : CG/JU
Sample : PB145686BL
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK686

Manual IntegrationsAPPROVED

Quant Time: Jun 21 23:08:14 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

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



TIC: BG053986.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.881min (+ 0.000) 33.61 ng/ul

response 22028

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.90	83.53
41.00	51.40	39.73#
42.00	37.10	27.86#

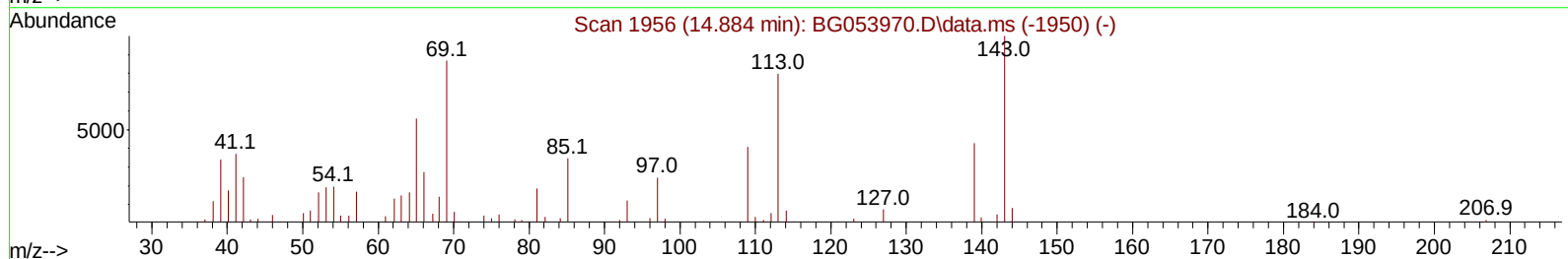
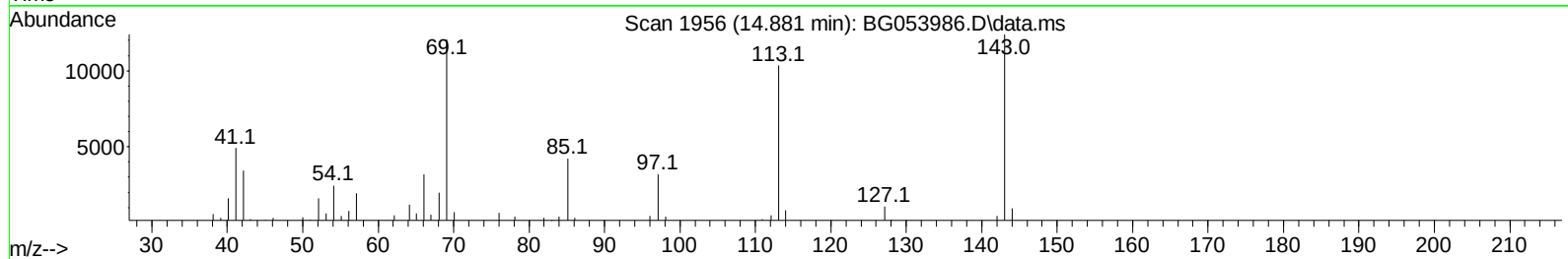
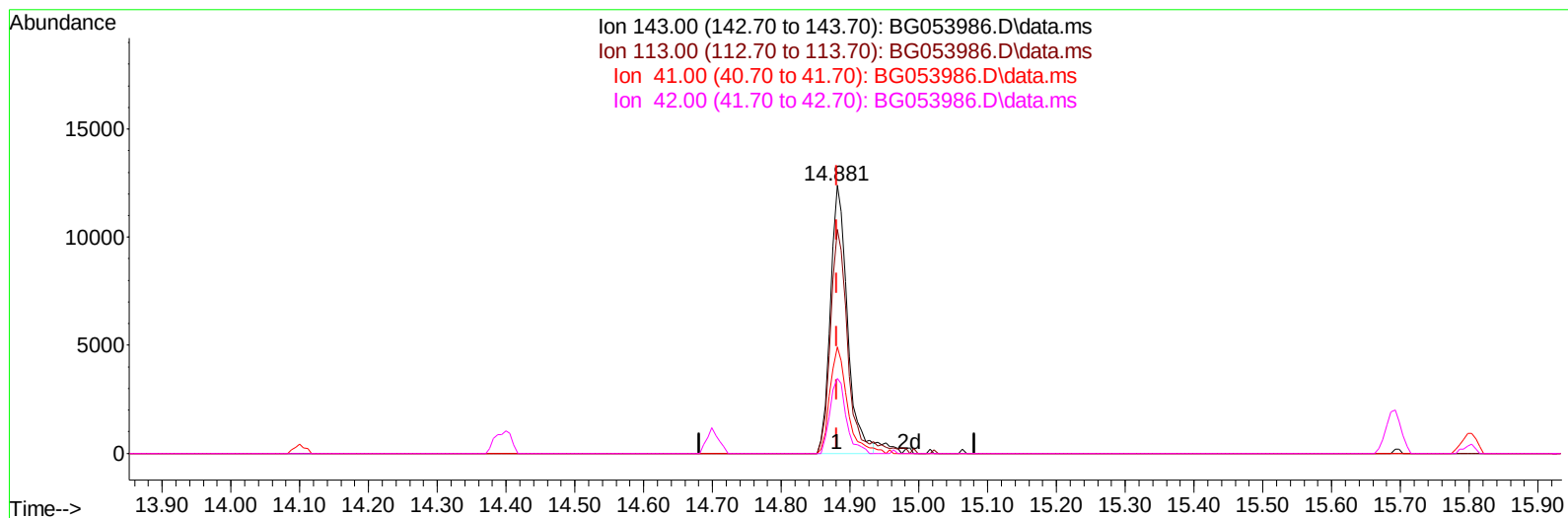
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TIC: BG053986.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.881min (+ 0.000) 32.39 ng/ul m

response 21232

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.90	83.53
41.00	51.40	39.73#
42.00	37.10	27.86#

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 Sample : PB145686BL
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.078	152	18747	20.000	ng/ul	0.00
20) Naphthalene-d8	10.886	136	77269	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.699	164	60846	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.443	188	166397	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.720	240	190755	20.000	ng/ul	# 0.00
88) Perylene-d12	24.975	264	191635	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.506	96	2632	6.680	ng/uL	0.00
4) Pyridine-d5	3.918	84	38831	36.234	ng/ul	0.00
7) Phenol-d5	7.231	99	49213	36.643	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.396	67	29611	36.549	ng/ul	0.00
11) 2-Chlorophenol-d4	7.607	132	39875	37.231	ng/ul	0.00
15) 4-Methylphenol-d8	8.777	113	41668	36.386	ng/ul	0.00
21) Nitrobenzene-d5	9.241	128	20246	35.300	ng/ul	0.00
24) 2-Nitrophenol-d4	9.964	143	23343	38.306	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.504	165	46607	34.569	ng/ul	0.00
31) 4-Chloroaniline-d4	11.015	131	58873	34.565	ng/ul	0.00
46) Dimethylphthalate-d6	14.100	166	173203	36.184	ng/ul	0.00
49) Acenaphthylene-d8	14.394	160	182986	35.156	ng/ul	0.00
54) 4-Nitrophenol-d4	14.881	143	21232m	32.394	ng/ul	0.00
60) Fluorene-d10	15.692	176	153069	35.601	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	27167	29.929	ng/ul	0.00
73) Anthracene-d10	17.543	188	270257	35.964	ng/ul	0.00
81) Pyrene-d10	19.828	212	345188	34.707	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.752	264	351123	36.298	ng/ul	0.00

Target Compounds Qvalue

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

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