

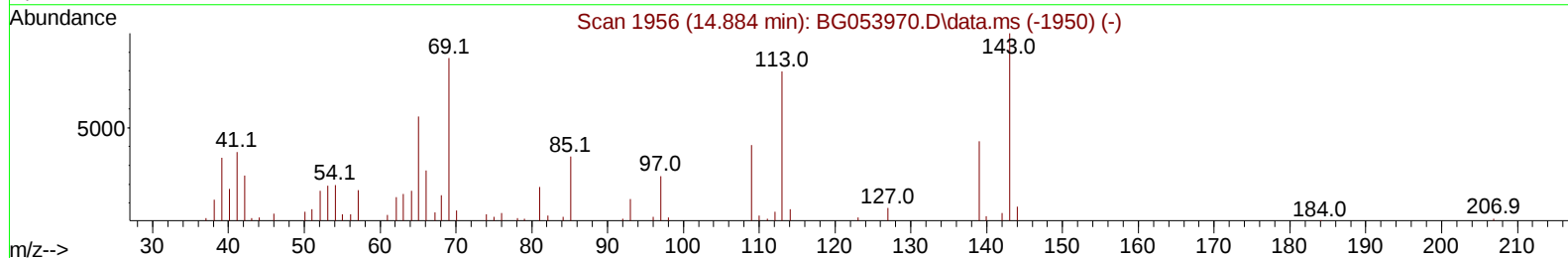
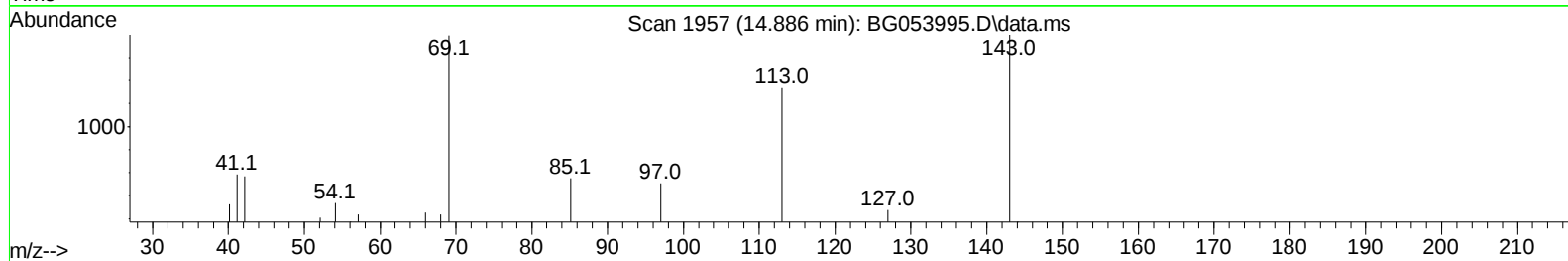
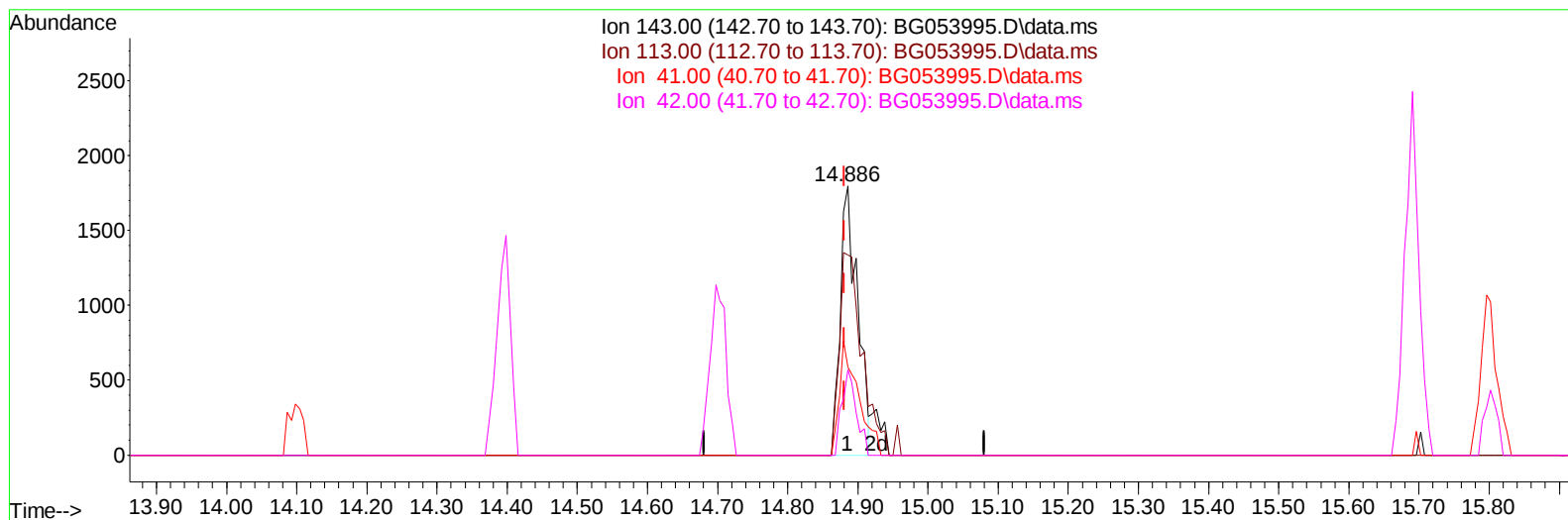
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062122\
 Data File : BG053995.D
 Acq On : 22 Jun 2022 2:30
 Operator : CG/JU
 Sample : N3355-19
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EY0D0

Manual IntegrationsAPPROVED

Quant Time: Jun 22 03:11:07 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: BG053995.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.886min (+ 0.005) 4.38 ng/u1

response 3073

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.90	74.39
41.00	51.40	32.57#
42.00	37.10	31.74

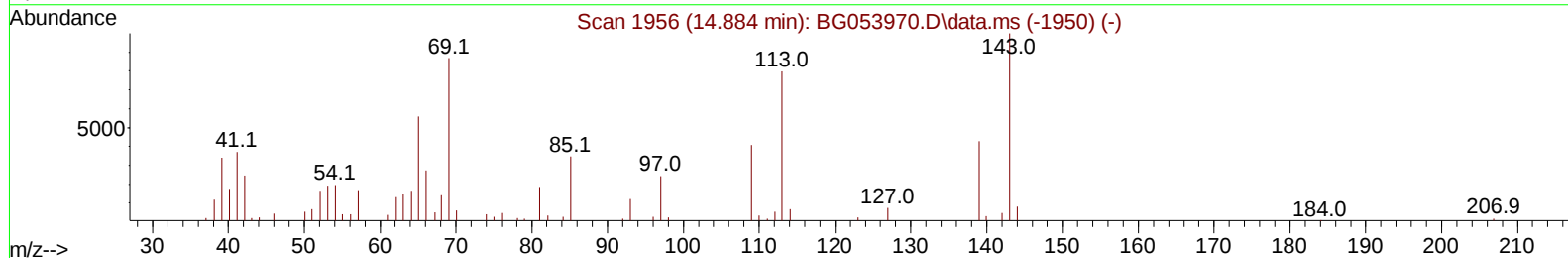
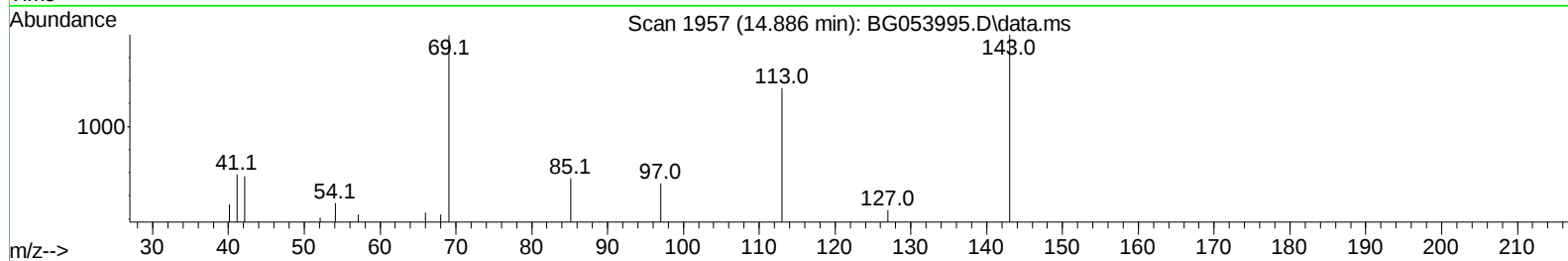
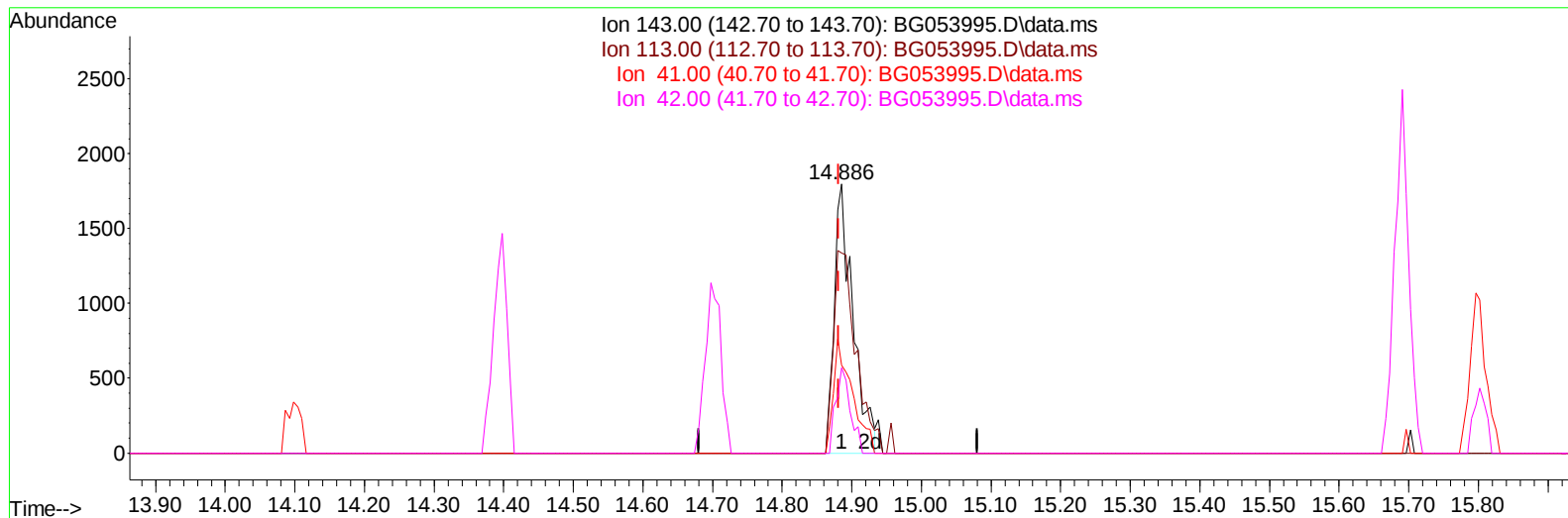
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TIC: BG053995.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.886min (+ 0.005) 4.75 ng/ul m

response 3335

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.90	74.39
41.00	51.40	32.57#
42.00	37.10	31.74

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.082	152	21166	20.000	ng/ul	0.00
20) Naphthalene-d8	10.884	136	84198	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.703	164	65112	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.447	188	178866	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.719	240	201534	20.000	ng/ul	# 0.00
88) Perylene-d12	24.980	264	201422	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.505	96	2606	5.858	ng/uL	0.00
4) Pyridine-d5	3.922	84	10335	8.542	ng/ul	0.00
7) Phenol-d5	7.236	99	9293	6.129	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.400	67	30629	33.485	ng/ul	0.00
11) 2-Chlorophenol-d4	7.612	132	30791	25.464	ng/ul	0.00
15) 4-Methylphenol-d8	8.775	113	20474	15.836	ng/ul	0.00
21) Nitrobenzene-d5	9.239	128	21960	35.138	ng/ul	0.00
24) 2-Nitrophenol-d4	9.962	143	23728	35.733	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.502	165	44193	30.081	ng/ul	0.00
31) 4-Chloroaniline-d4	11.014	131	52607	28.344	ng/ul	0.00
46) Dimethylphthalate-d6	14.104	166	184064	35.934	ng/ul	0.00
49) Acenaphthylene-d8	14.398	160	196673	35.310	ng/ul	0.00
54) 4-Nitrophenol-d4	14.886	143	3335m	4.755	ng/ul	0.00
60) Fluorene-d10	15.691	176	167657	36.439	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.802	200	29292	30.021	ng/ul	0.00
73) Anthracene-d10	17.541	188	301944	37.379	ng/ul	0.00
81) Pyrene-d10	19.827	212	390978	37.208	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.750	264	392082	38.563	ng/ul	0.00

Target Compounds Qvalue

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

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