

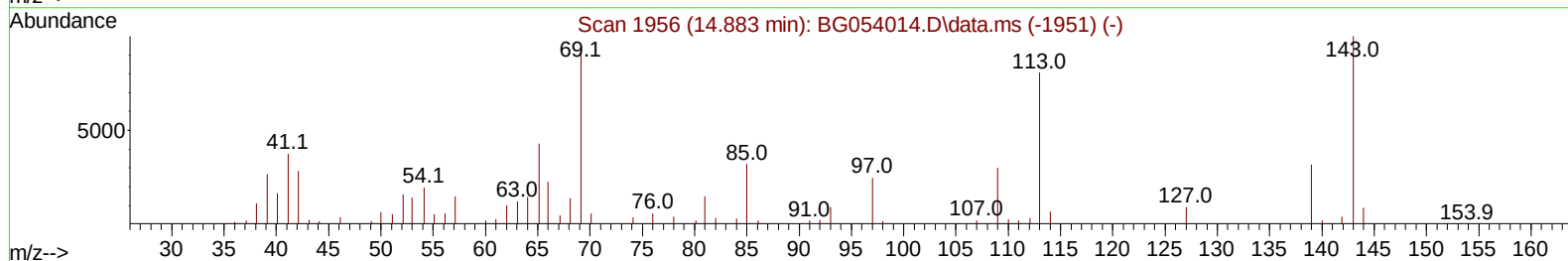
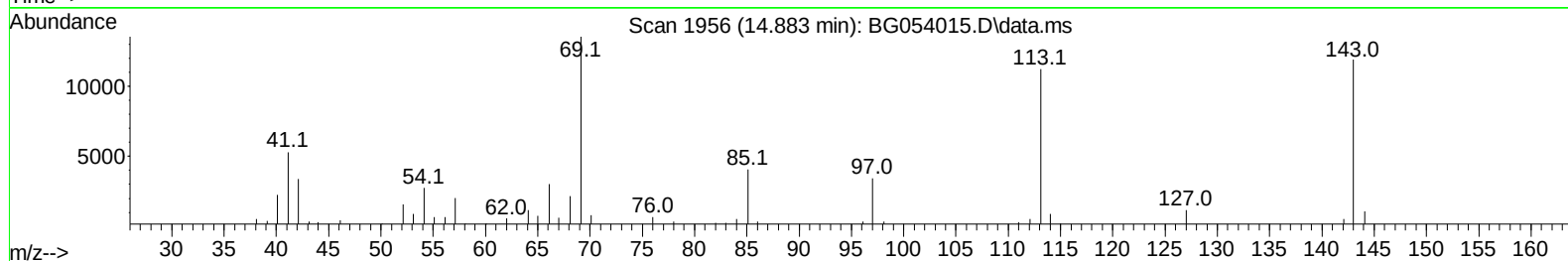
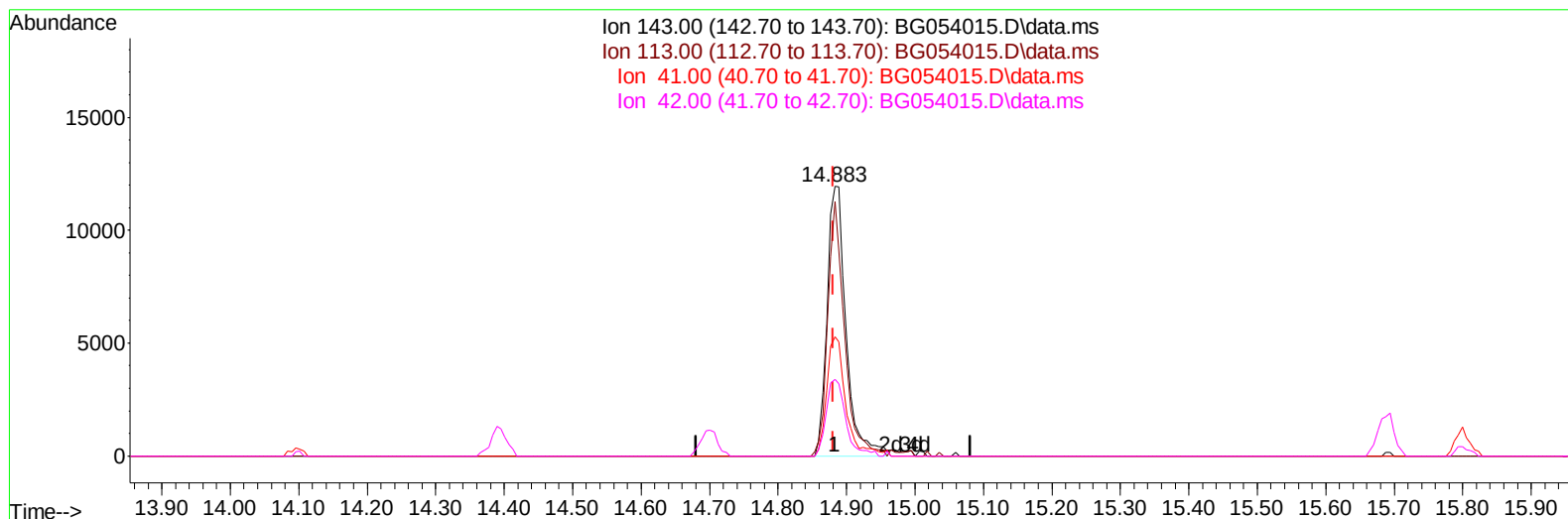
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062322\
Data File : BG054015.D
Acq On : 23 Jun 2022 16:20
Operator : CG/JU
Sample : PB145673BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK673

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 06/24/2022
Supervised By :mohammad ahmed 06/29/2022

Quant Time: Jun 23 23:05:01 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: BG054015.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.883min (+ 0.002) 32.32 ng/ul

response 22973

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.90	94.34
41.00	51.40	44.34
42.00	37.10	28.34#

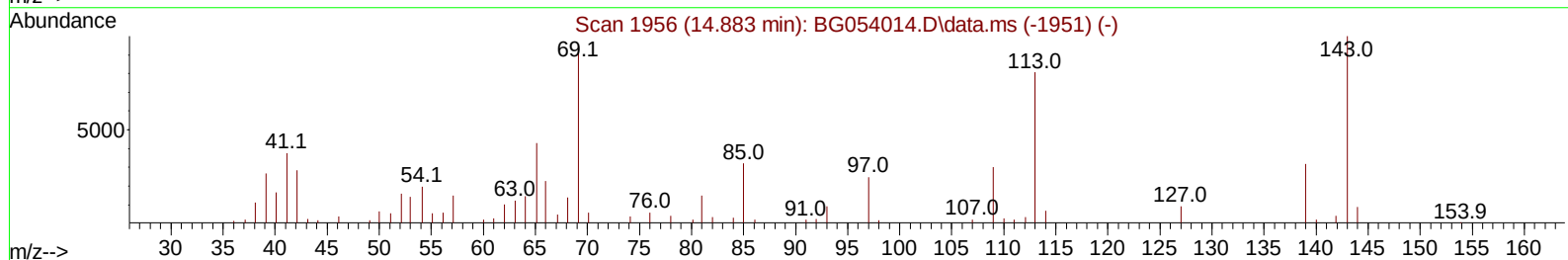
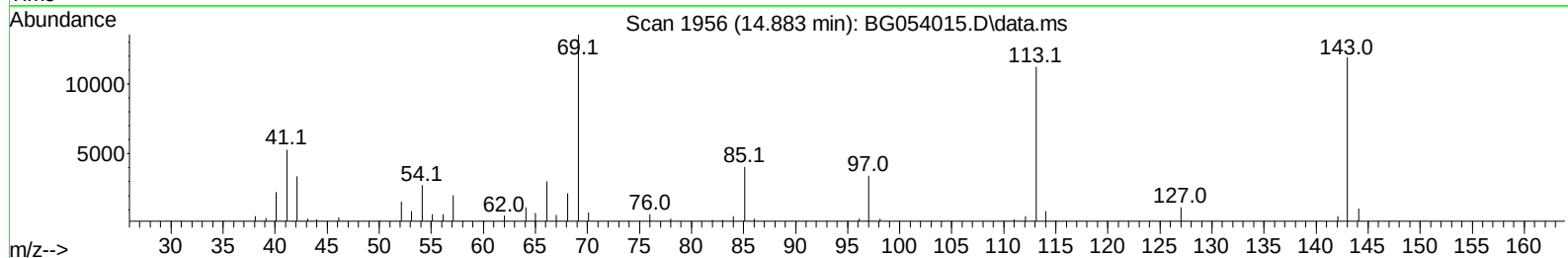
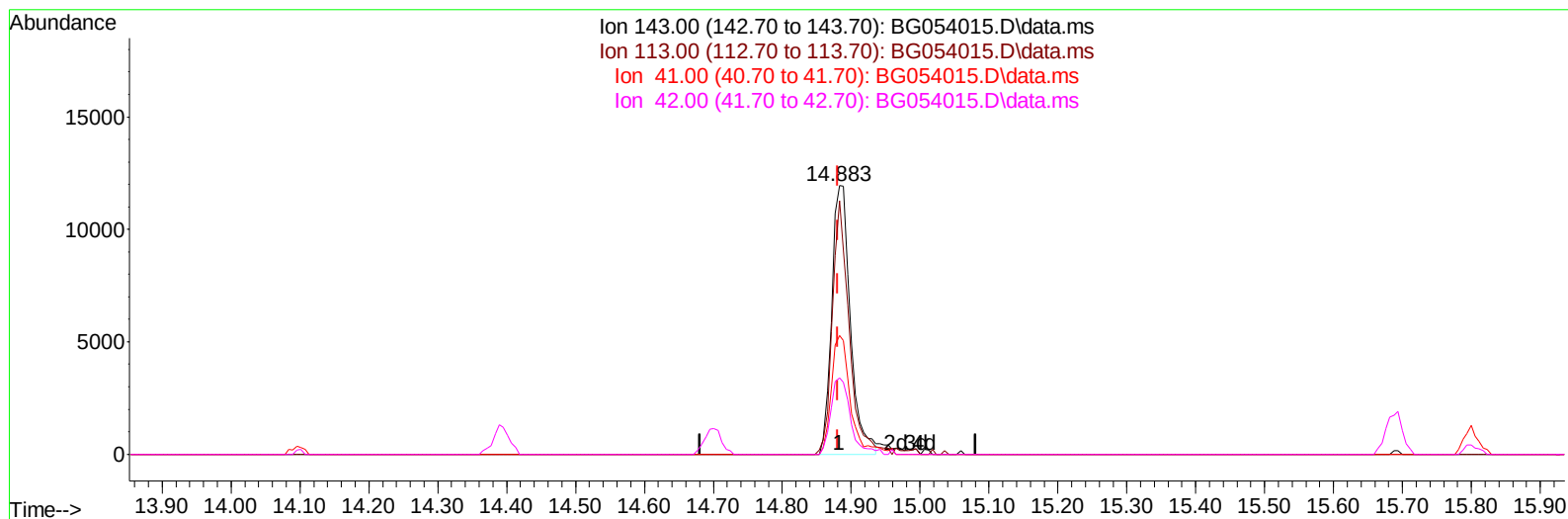
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TIC: BG054015.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.883min (+ 0.002) 31.57 ng/ul m

response 22443

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.90	94.34
41.00	51.40	44.34
42.00	37.10	28.34#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.073	152	17856	20.000	ng/ul	0.00
20) Naphthalene-d8	10.888	136	80267	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.701	164	65989	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.445	188	170181	20.000	ng/ul	0.00
79) Chrysene-d12	21.716	240	194411	20.000	ng/ul	0.00
88) Perylene-d12	24.983	264	191465	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.496	96	2362	6.294	ng/uL	0.00
4) Pyridine-d5	3.908	84	35789	35.062	ng/ul	0.00
7) Phenol-d5	7.227	99	47682	37.274	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.392	67	27868	36.115	ng/ul	0.00
11) 2-Chlorophenol-d4	7.603	132	37479	36.740	ng/ul	0.00
15) 4-Methylphenol-d8	8.772	113	41723	38.253	ng/ul	0.00
21) Nitrobenzene-d5	9.231	128	19666	33.008	ng/ul	0.00
24) 2-Nitrophenol-d4	9.959	143	22519	35.573	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.500	165	47449	33.879	ng/ul	0.00
31) 4-Chloroaniline-d4	11.011	131	62036	35.061	ng/ul	0.00
46) Dimethylphthalate-d6	14.101	166	176939	34.084	ng/ul	0.00
49) Acenaphthylene-d8	14.395	160	190885	33.815	ng/ul	0.00
54) 4-Nitrophenol-d4	14.883	143	22443m	31.573	ng/ul	0.00
60) Fluorene-d10	15.688	176	157981	33.879	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.799	200	27222	29.323	ng/ul	0.00
73) Anthracene-d10	17.539	188	267558	34.813	ng/ul	0.00
81) Pyrene-d10	19.824	212	358497	35.367	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.754	264	342301	35.418	ng/ul	0.00

Target Compounds Qvalue

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

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