

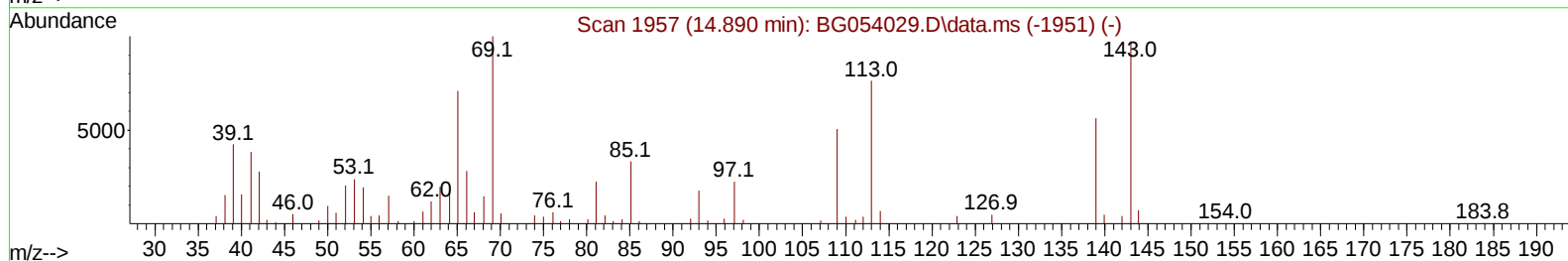
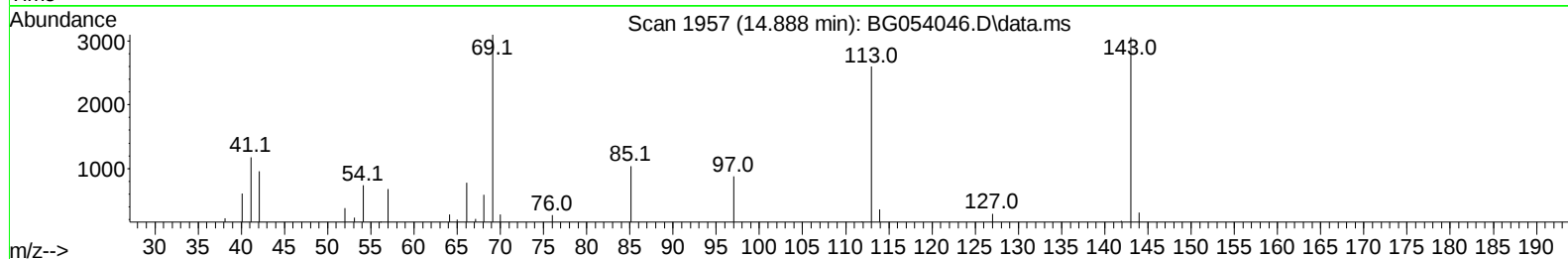
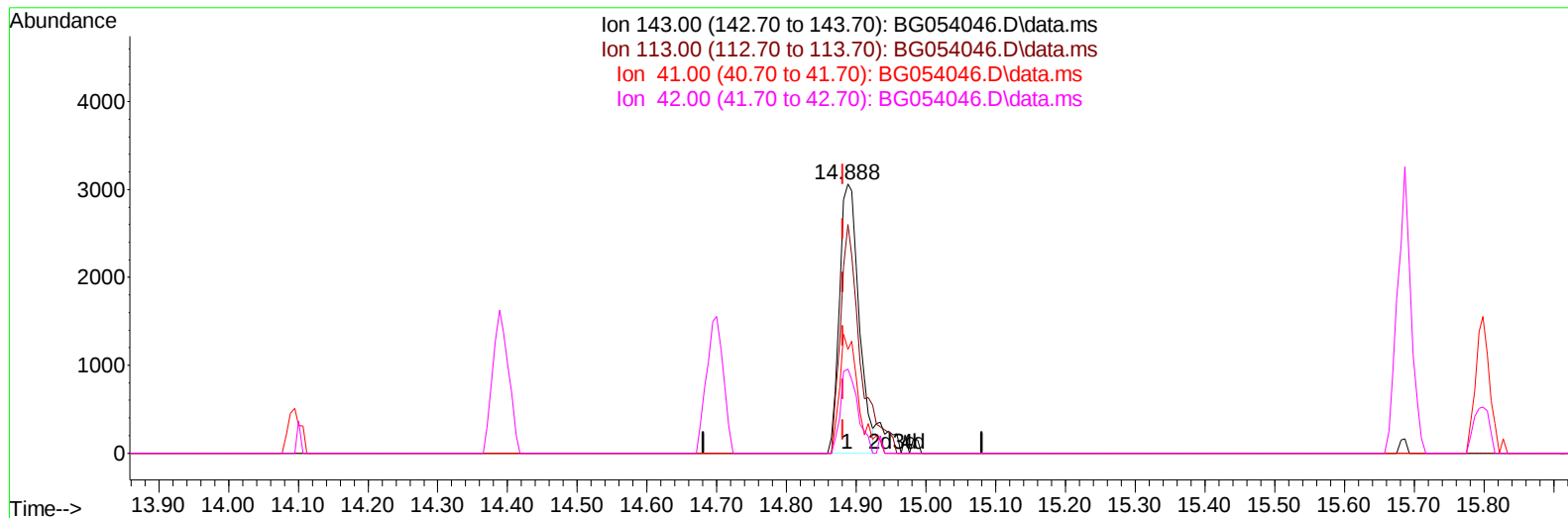
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062322\
Data File : BG054046.D
Acq On : 24 Jun 2022 13:52
Operator : CG/JU
Sample : N3395-19
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0166

Manual IntegrationsAPPROVED

Quant Time: Jun 24 23:53:11 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 06/27/2022
Supervised By :mohammad ahmed 06/29/2022



TIC: BG054046.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.888min (+ 0.007) 5.57 ng/u1

response 5878

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.90	84.91
41.00	51.40	38.52#
42.00	37.10	31.20

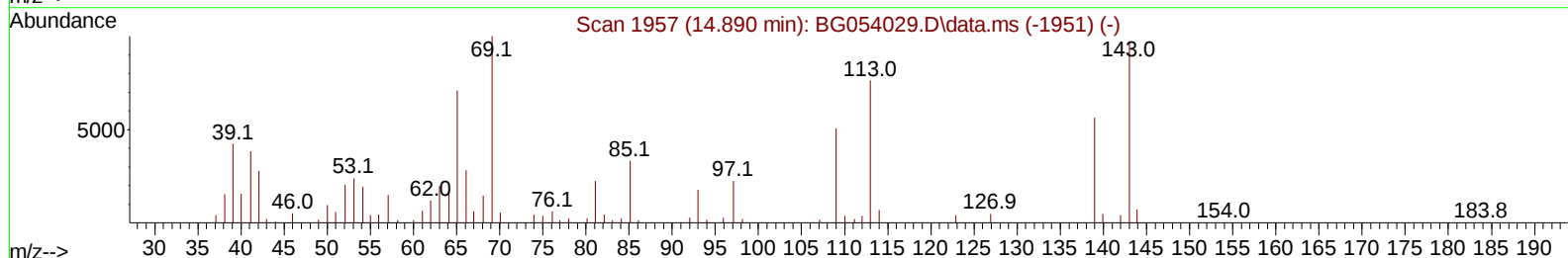
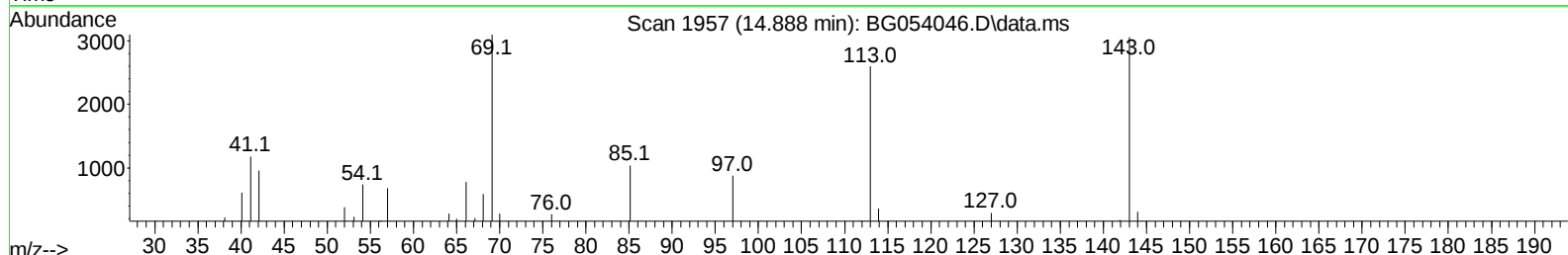
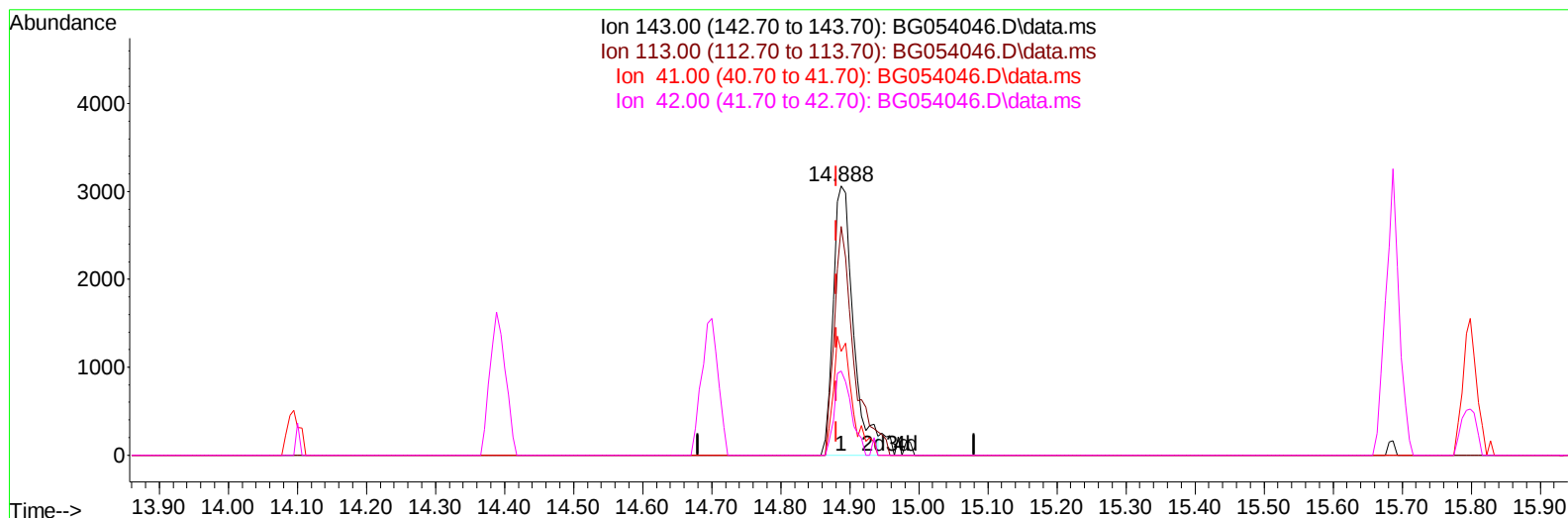
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TIC: BG054046.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.888min (+ 0.007) 5.87 ng/ul m

response 6198

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	83.90	84.91
41.00	51.40	38.52#
42.00	37.10	31.20

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.072	152	24005	20.000	ng/ul	0.00
20) Naphthalene-d8	10.881	136	116218	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.700	164	97940	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.438	188	250199	20.000	ng/ul	0.00
79) Chrysene-d12	21.715	240	278008	20.000	ng/ul	0.00
88) Perylene-d12	24.976	264	270911	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.495	96	2023	4.010	ng/uL	-0.01
4) Pyridine-d5	3.918	84	8610	6.274	ng/ul	0.00
7) Phenol-d5	7.232	99	14230	8.274	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.391	67	35167	33.900	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.602	132	36662	26.733	ng/ul	0.00
15) 4-Methylphenol-d8	8.777	113	28283	19.288	ng/ul	0.00
21) Nitrobenzene-d5	9.230	128	24175	28.025	ng/ul	0.00
24) 2-Nitrophenol-d4	9.958	143	27039	29.501	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.499	165	54477	26.864	ng/ul	0.00
31) 4-Chloroaniline-d4	11.010	131	57411	22.410	ng/ul	0.00
46) Dimethylphthalate-d6	14.095	166	227888	29.577	ng/ul	0.00
49) Acenaphthylene-d8	14.394	160	245902	29.350	ng/ul	0.00
54) 4-Nitrophenol-d4	14.888	143	6198m	5.875	ng/ul	0.00
60) Fluorene-d10	15.687	176	200229	28.931	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.799	200	35023	25.661	ng/ul	0.00
73) Anthracene-d10	17.538	188	353379	31.274	ng/ul	0.00
81) Pyrene-d10	19.823	212	451750	31.166	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.753	264	440823	32.236	ng/ul	0.00

Target Compounds Qvalue

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

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