

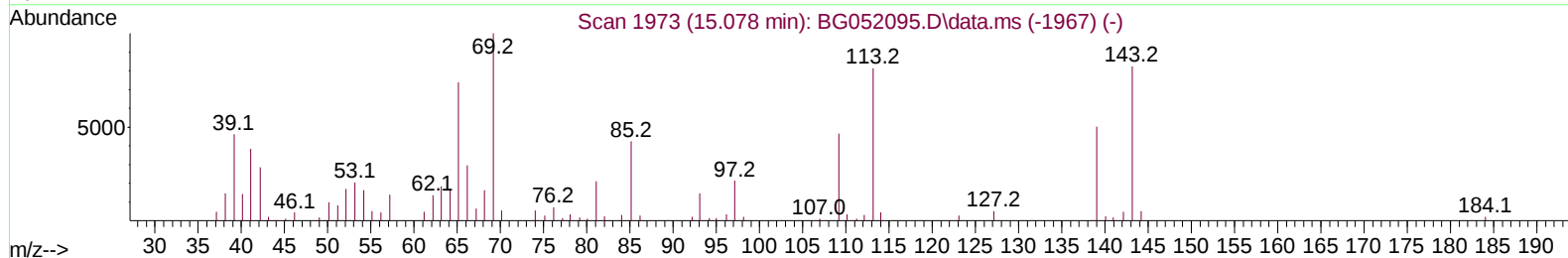
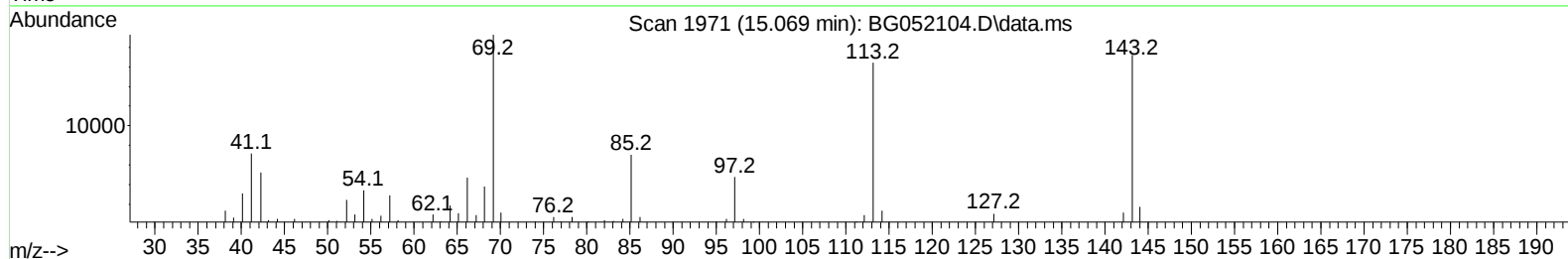
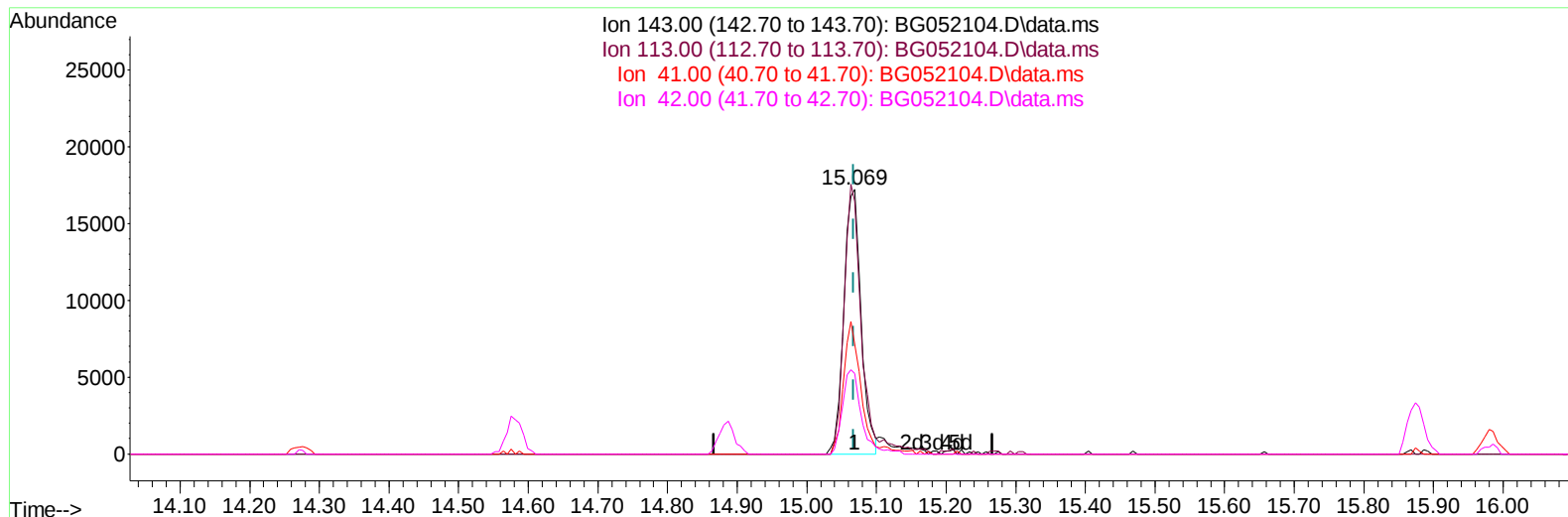
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011922\
Data File : BG052104.D
Acq On : 20 Jan 2022 22:32
Operator : CG/JU
Sample : PB142157BL
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK157

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/21/2022
Supervised By :Yogesh Patel 01/23/2022

Quant Time: Jan 21 00:01:49 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jan 06 14:43:02 2022
Response via : Initial Calibration



TIC: BG052104.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.069min (+ 0.002) 22.34 ng/u1

response 29756

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	95.45
41.00	44.40	41.43
42.00	29.70	30.40

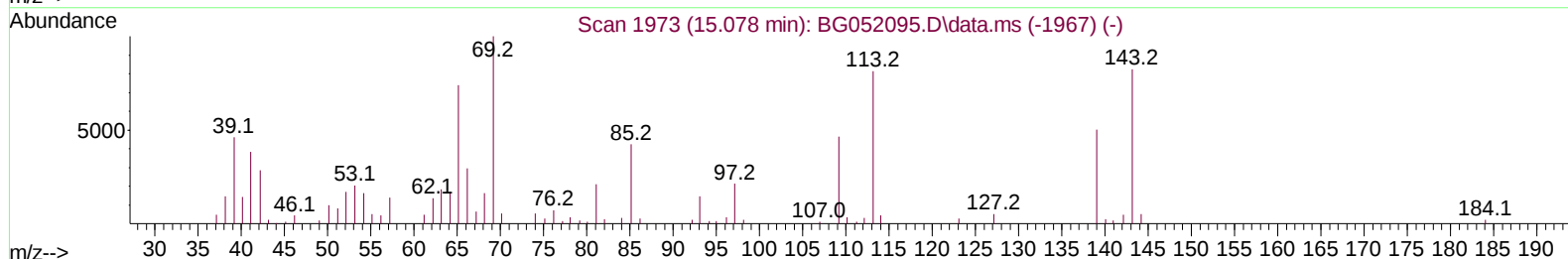
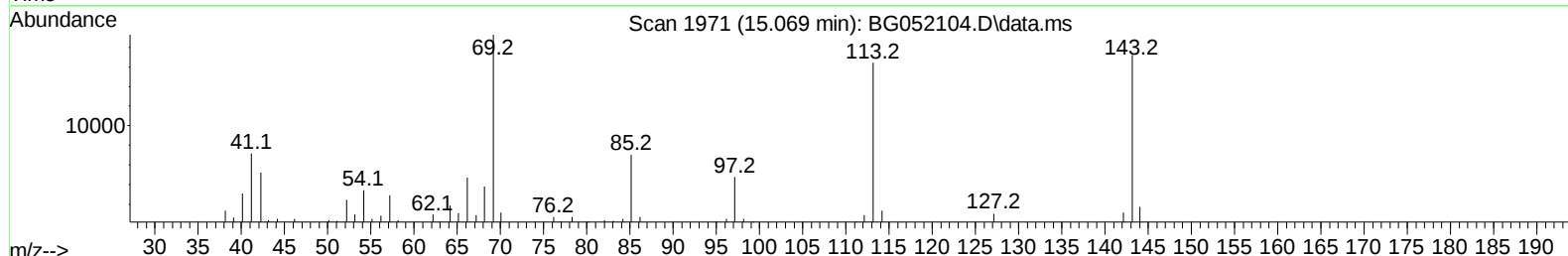
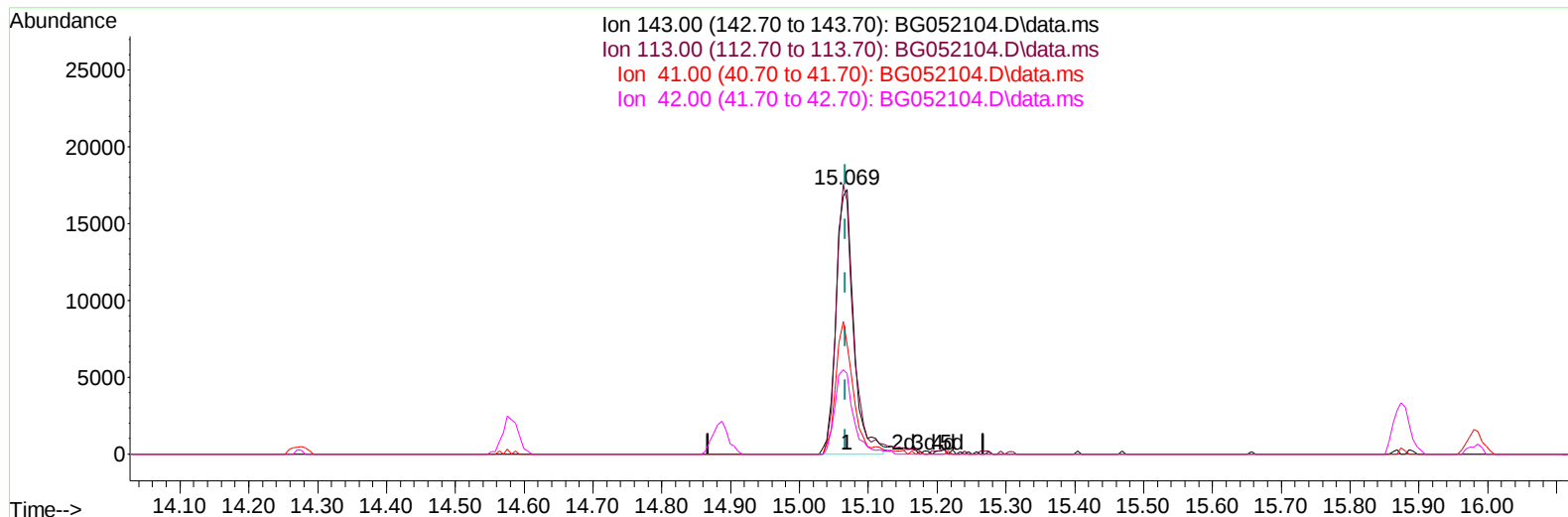
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TIC: BG052104.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.069min (+ 0.002) 23.19 ng/ul m

response 30876

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	95.45
41.00	44.40	41.43
42.00	29.70	30.40

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.243	152	29848	20.000	ng/ul	0.00
20) Naphthalene-d8	11.080	136	130066	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.887	164	96364	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.636	188	220004	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.954	240	212724	20.000	ng/ul	# 0.00
88) Perylene-d12	25.438	264	211862	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.567	96	3964	4.746	ng/uL	0.00
4) Pyridine-d5	3.996	84	58348	22.471	ng/ul	0.00
7) Phenol-d5	7.380	99	76367	26.141	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.550	67	45762	24.677	ng/ul	0.00
11) 2-Chlorophenol-d4	7.773	132	54947	28.409	ng/ul	0.00
15) 4-Methylphenol-d8	8.942	113	64224	28.281	ng/ul	0.00
21) Nitrobenzene-d5	9.412	128	30104	29.291	ng/ul	0.00
24) 2-Nitrophenol-d4	10.146	143	34894	30.699	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.693	165	66906	30.114	ng/ul	0.00
31) 4-Chloroaniline-d4	11.204	131	83810	27.087	ng/ul	0.00
46) Dimethylphthalate-d6	14.276	166	224753	28.217	ng/ul	0.00
49) Acenaphthylene-d8	14.582	160	275213	28.604	ng/ul	0.00
54) 4-Nitrophenol-d4	15.069	143	30876m	23.185	ng/ul	0.00
60) Fluorene-d10	15.874	176	192630	27.738	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.980	200	30505	21.979	ng/ul	0.00
73) Anthracene-d10	17.736	188	315513	30.232	ng/ul	0.00
81) Pyrene-d10	20.010	212	376603	30.762	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.197	264	354015	31.741	ng/ul	0.00

Target Compounds Qvalue

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

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