

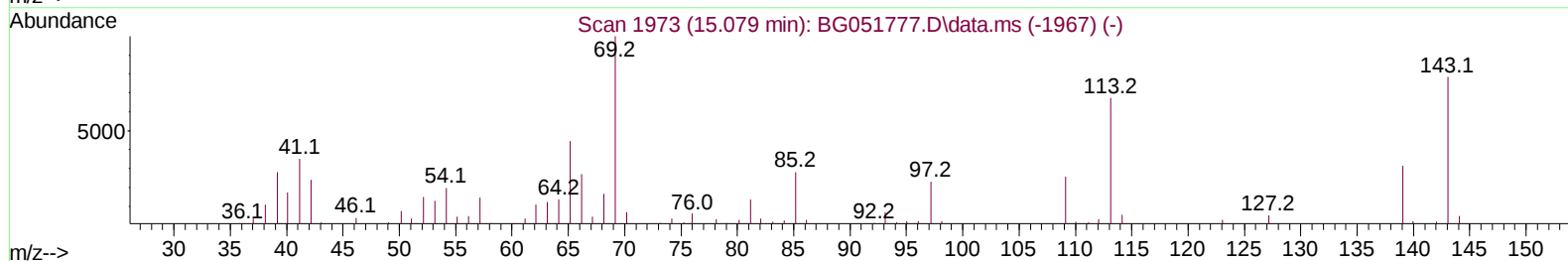
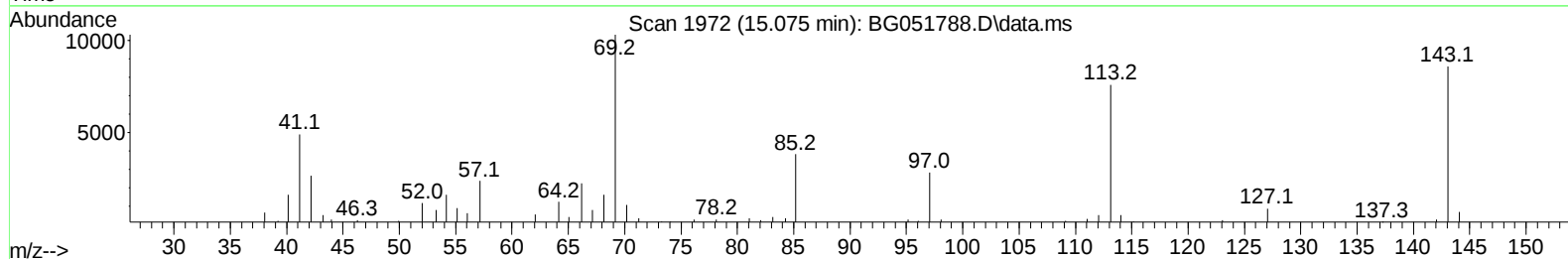
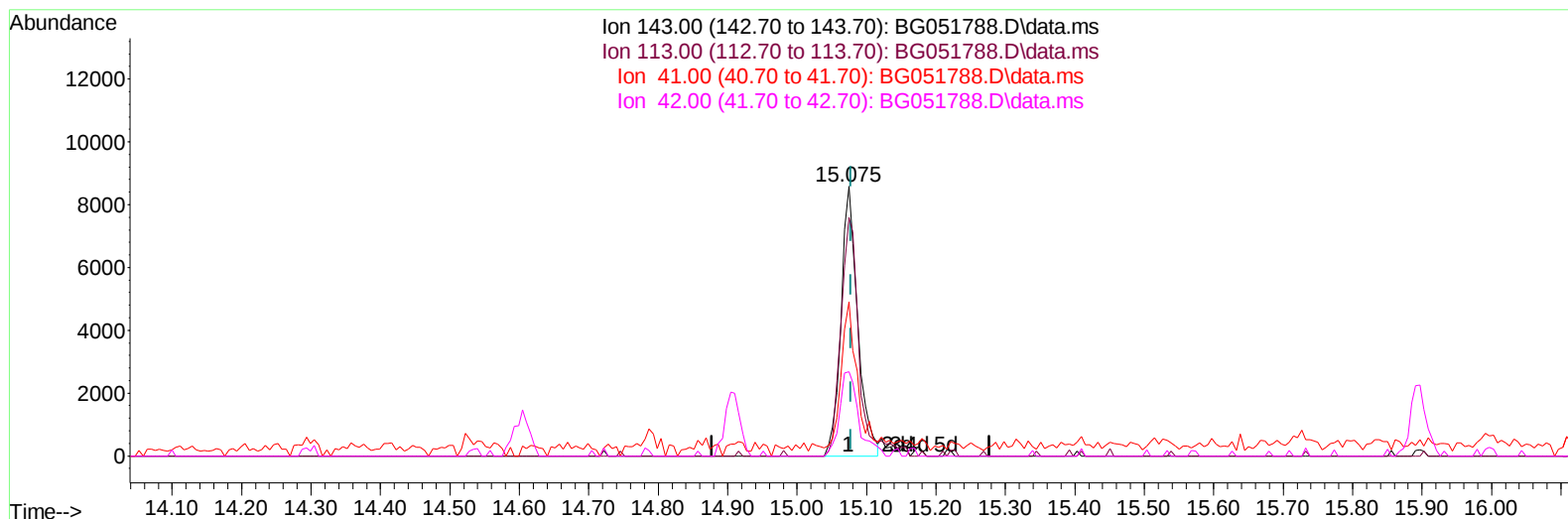
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051788.D
Acq On : 23 Dec 2021 17:04
Operator : CG/JU
Sample : M5215-03
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD5

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/27/2021
Supervised By :mohammad ahmed 12/31/2021

Quant Time: Dec 24 00:12:18 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 23 13:37:07 2021
Response via : Initial Calibration



TIC: BG051788.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.075min (-0.003) 12.40 ng/u1

response 14319

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	88.47
41.00	44.40	57.19#
42.00	29.70	31.25

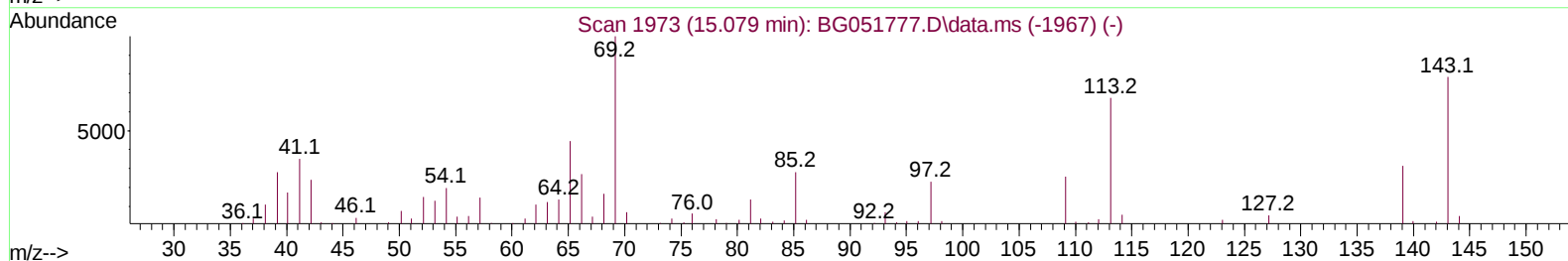
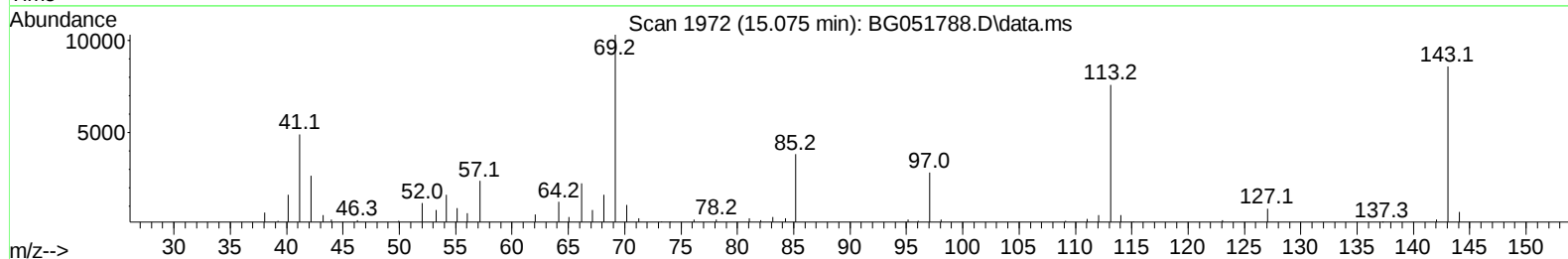
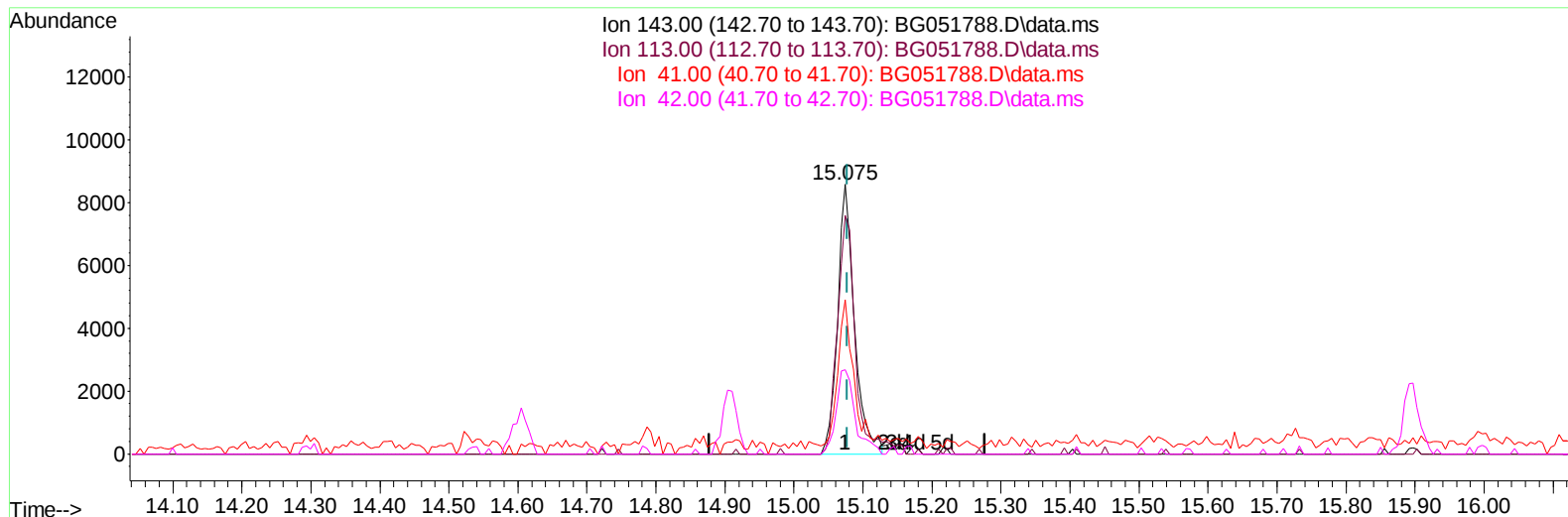
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(54) 4-Nitrophenol-d4 (S)

15.075min (-0.003) 12.65 ng/ul m

response 14607

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	88.47
41.00	44.40	57.19#
42.00	29.70	31.25

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.272	152	31110	20.000	ng/ul	-0.02
20) Naphthalene-d8	11.109	136	124300	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.910	164	88733	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.653	188	184525	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.977	240	147314	20.000	ng/ul	0.00
88) Perylene-d12	25.472	264	157267	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.596	96	2198	2.934	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.403	99	55759	20.074	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.579	67	33792	20.464	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.796	132	38472	19.914	ng/ul	-0.01
15) 4-Methylphenol-d8	8.971	113	45125	21.017	ng/ul	0.00
21) Nitrobenzene-d5	9.441	128	21371	23.451	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.169	143	22175	22.259	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.716	165	43585	20.175	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.233	131	47938	16.853	ng/ul	0.00
46) Dimethylphthalate-d6	14.293	166	142485	20.062	ng/ul	-0.02
49) Acenaphthylene-d8	14.605	160	183072	20.755	ng/ul	0.00
54) 4-Nitrophenol-d4	15.075	143	14607m	12.647	ng/ul	0.00
60) Fluorene-d10	15.897	176	126183	19.831	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.003	200	6563	6.827	ng/ul	0.00
73) Anthracene-d10	17.753	188	191351	21.739	ng/ul	0.00
81) Pyrene-d10	20.032	212	192982	21.766	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.231	264	174543	21.093	ng/ul	-0.02
Target Compounds						
					Qvalue	
8) Phenol	7.432	94	5691	2.060	ng/ul#	70
30) Naphthalene	11.156	128	7760	1.159	ng/ul#	96
78) Di-n-butylphthalate	18.593	149	10470	1.013	ng/ul#	95
86) Bis(2-ethylhexyl)phtha...	21.842	149	5663	1.197	ng/ul#	96

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

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