

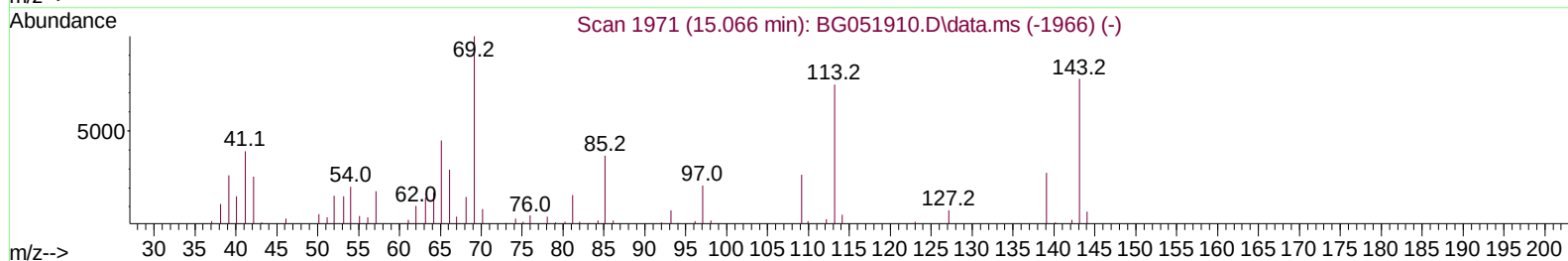
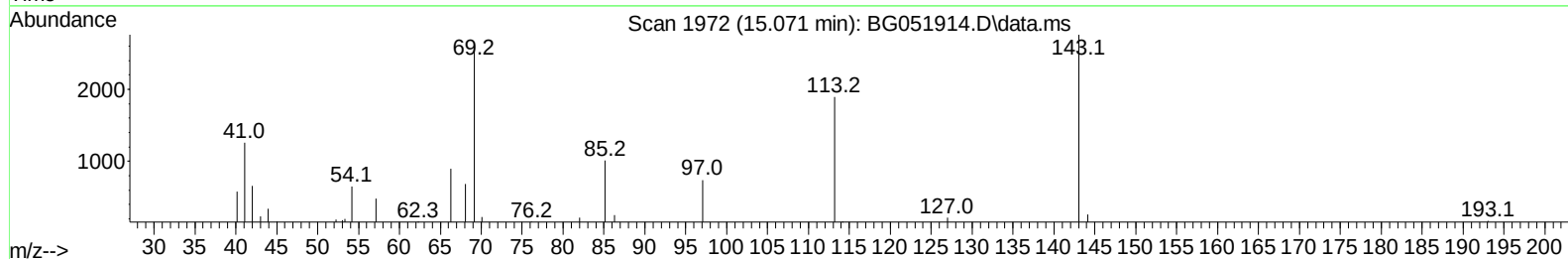
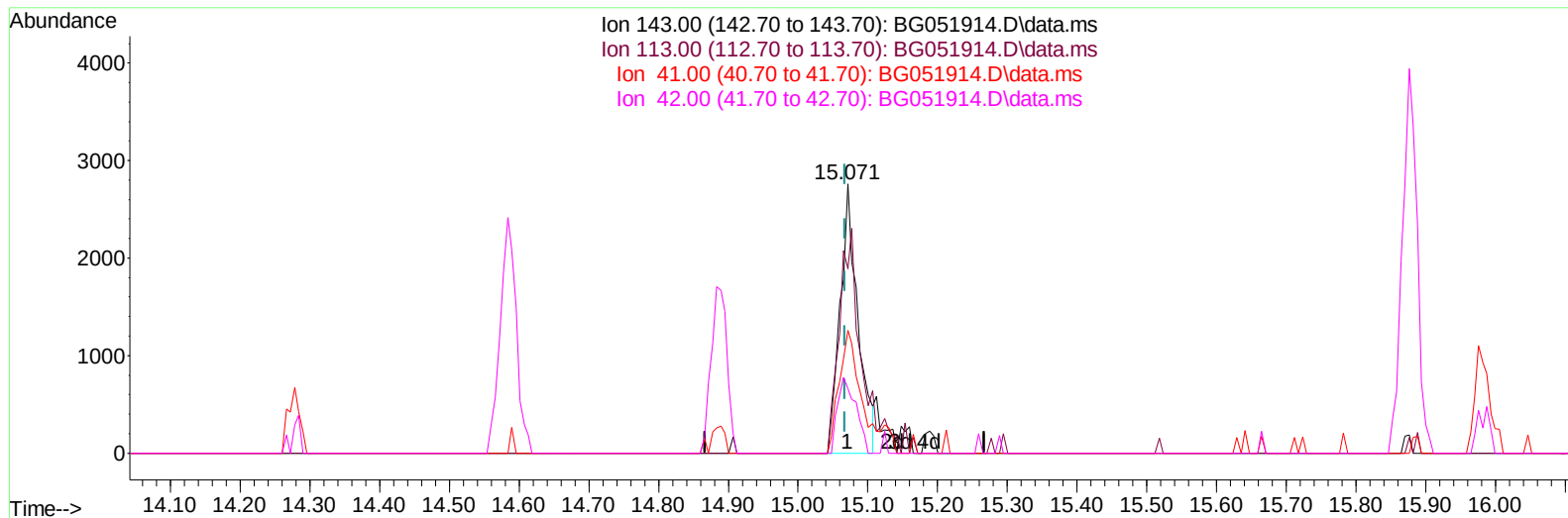
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010622\
 Data File : BG051914.D
 Acq On : 7 Jan 2022 14:16
 Operator : CG/JU
 Sample : N1027-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0J36

Manual IntegrationsAPPROVED

Quant Time: Jan 07 22:53:29 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

Reviewed By :Jagrut Upadhyay 01/10/2022
 Supervised By :mohammad ahmed 01/12/2022



TIC: BG051914.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.071min (+ 0.004) 5.26 ng/ul

response 4874

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	68.56
41.00	44.40	45.69
42.00	29.70	23.86

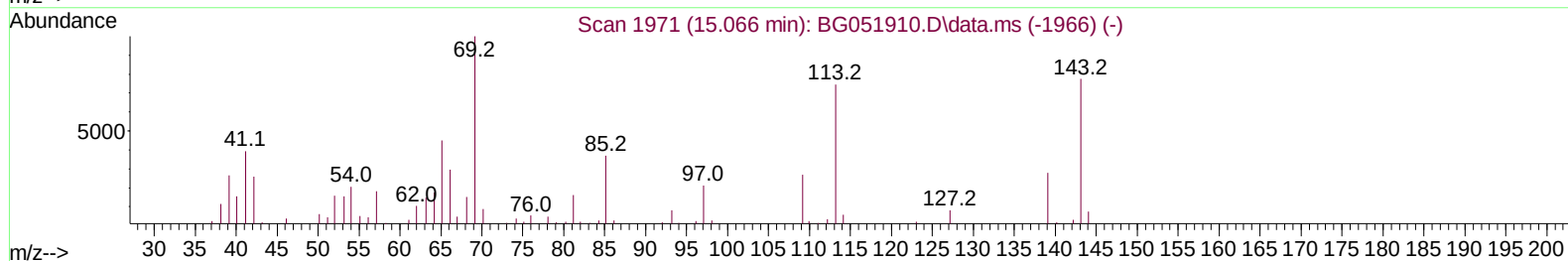
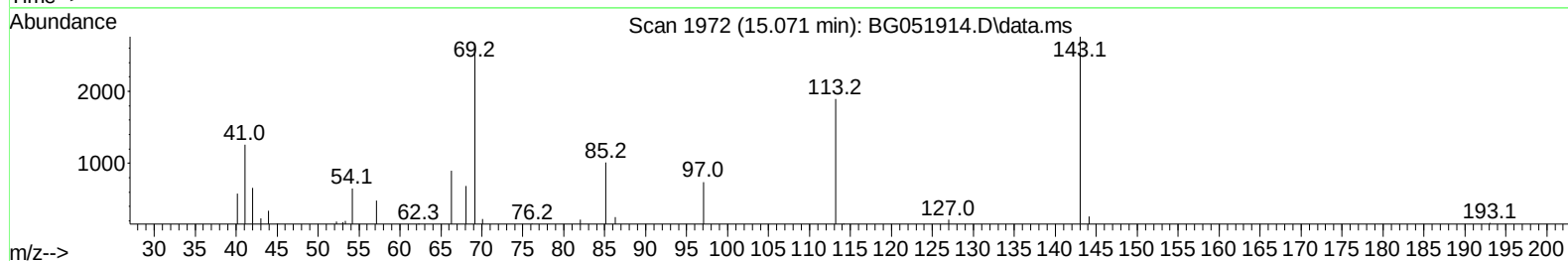
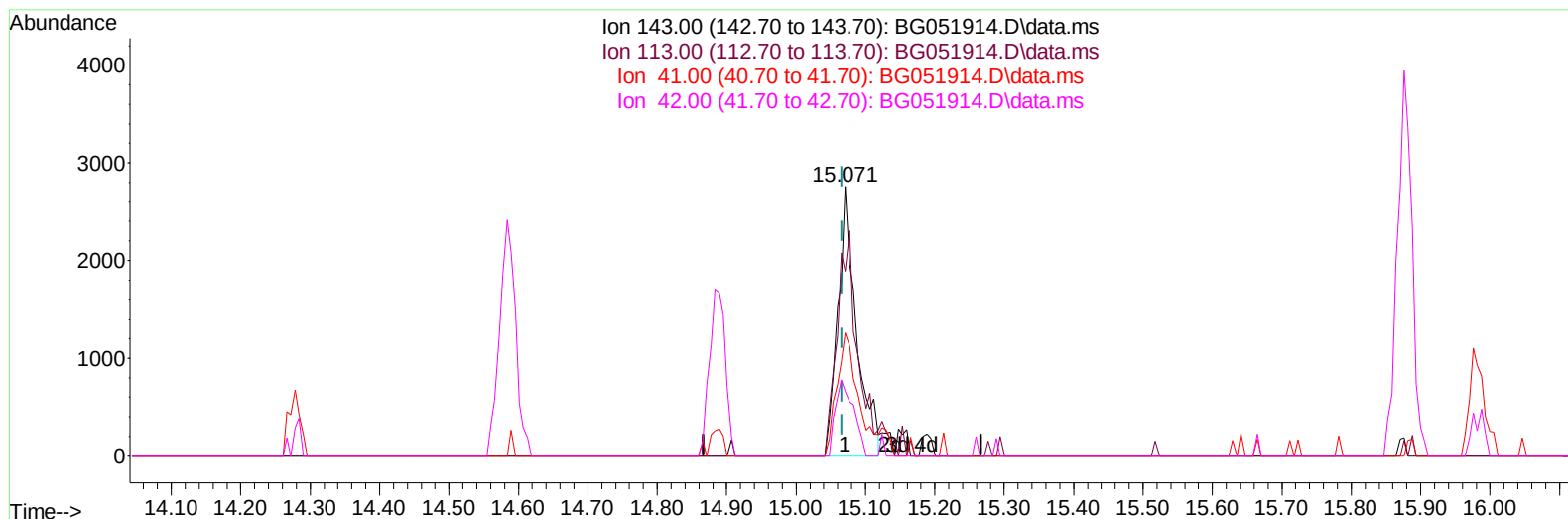
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(54) 4-Nitrophenol-d4 (S)

15.071min (+ 0.004) 5.56 ng/ul m

response 5157

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	68.56
41.00	44.40	45.69
42.00	29.70	23.86

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.245	152	20970	20.000	ng/ul	0.00
20) Naphthalene-d8	11.088	136	89516	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.889	164	67064	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.638	188	178303	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.956	240	192017	20.000	ng/ul	# 0.00
88) Perylene-d12	25.439	264	190901	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.575	96	2662	4.536	ng/uL	0.00
4) Pyridine-d5	4.010	84	16358	8.967	ng/ul	0.00
7) Phenol-d5	7.387	99	16177	7.882	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.552	67	47037	36.102	ng/ul	0.00
11) 2-Chlorophenol-d4	7.775	132	39382	28.982	ng/ul	0.00
15) 4-Methylphenol-d8	8.950	113	28318	17.749	ng/ul	0.00
21) Nitrobenzene-d5	9.414	128	26643	37.667	ng/ul	0.00
24) 2-Nitrophenol-d4	10.148	143	27000	34.514	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.695	165	52833	34.552	ng/ul	0.00
31) 4-Chloroaniline-d4	11.212	131	70378	33.050	ng/ul	0.00
46) Dimethylphthalate-d6	14.278	166	213572	38.528	ng/ul	0.00
49) Acenaphthylene-d8	14.584	160	248819	37.159	ng/ul	0.00
54) 4-Nitrophenol-d4	15.071	143	5157m	5.564	ng/ul	0.00
60) Fluorene-d10	15.882	176	185795	38.442	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.982	200	17321	15.399	ng/ul	0.00
73) Anthracene-d10	17.738	188	358449	42.378	ng/ul	0.00
81) Pyrene-d10	20.017	212	458403	41.481	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.205	264	420393	41.831	ng/ul	0.00

Target Compounds Qvalue

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

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