

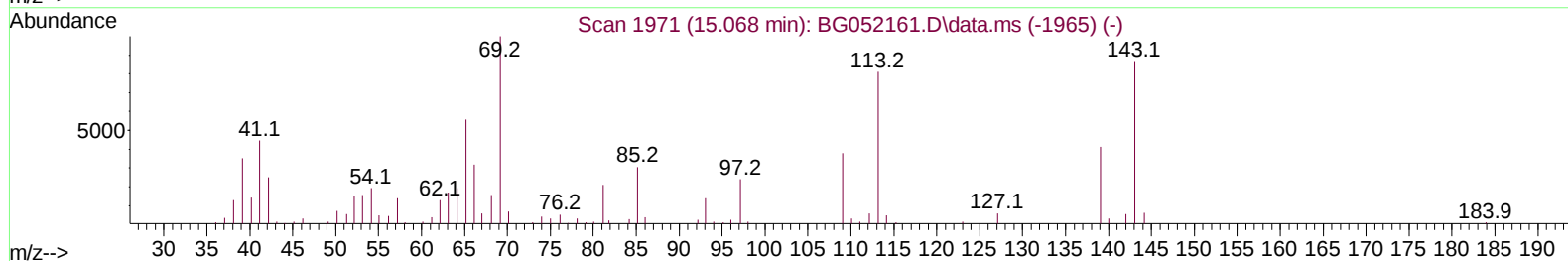
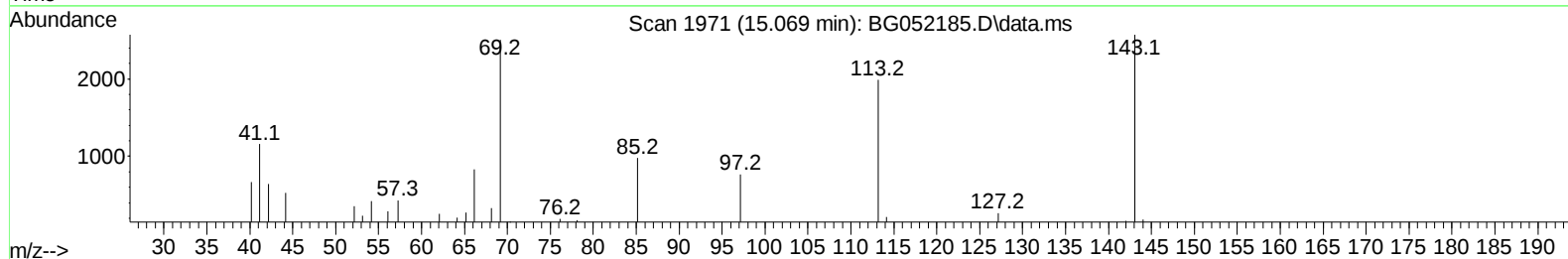
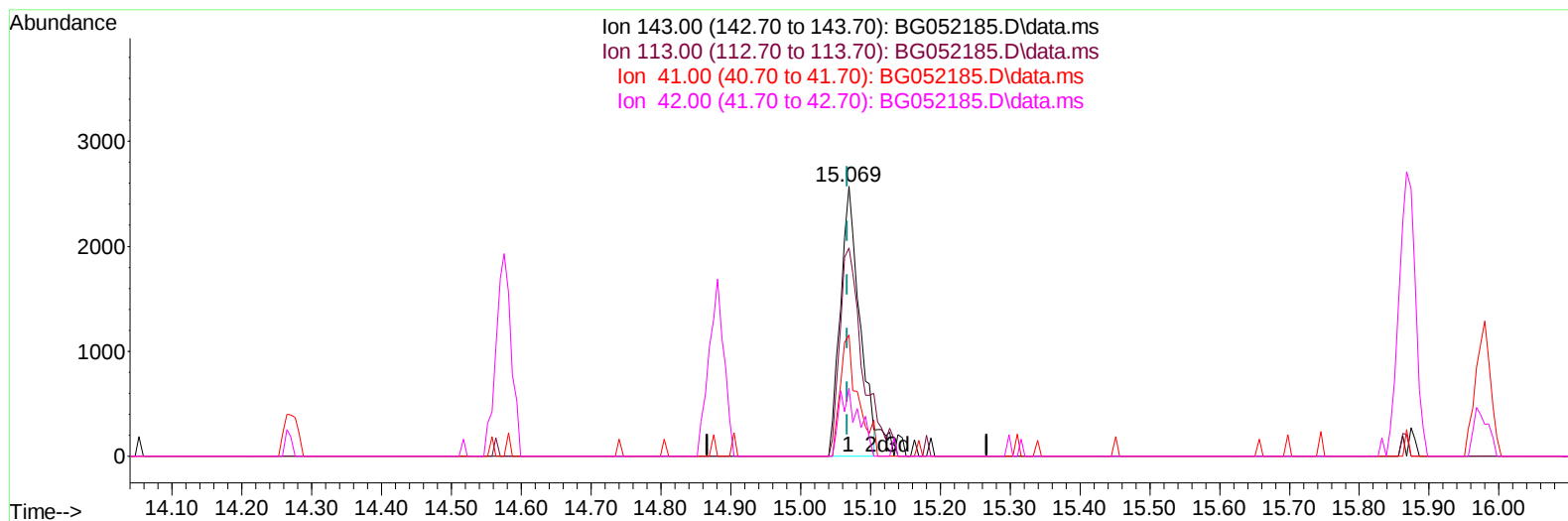
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052185.D
 Acq On : 27 Jan 2022 14:50
 Operator : CG/JU
 Sample : N1279-05
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0VD3

Manual IntegrationsAPPROVED

Quant Time: Jan 27 23:59:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 26 23:58:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/28/2022
 Supervised By :Yogesh Patel 01/28/2022



TIC: BG052185.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.069min (+ 0.002) 4.90 ng/u1

response 4866

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	77.20
41.00	44.40	45.02
42.00	29.70	25.10

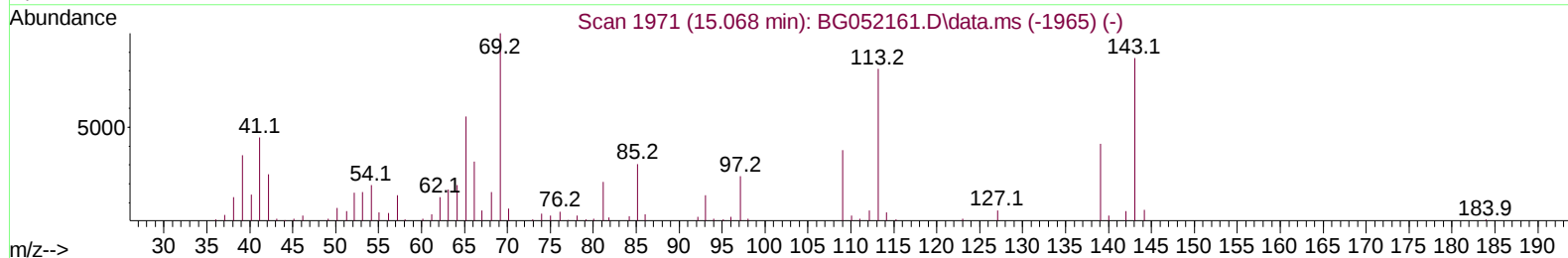
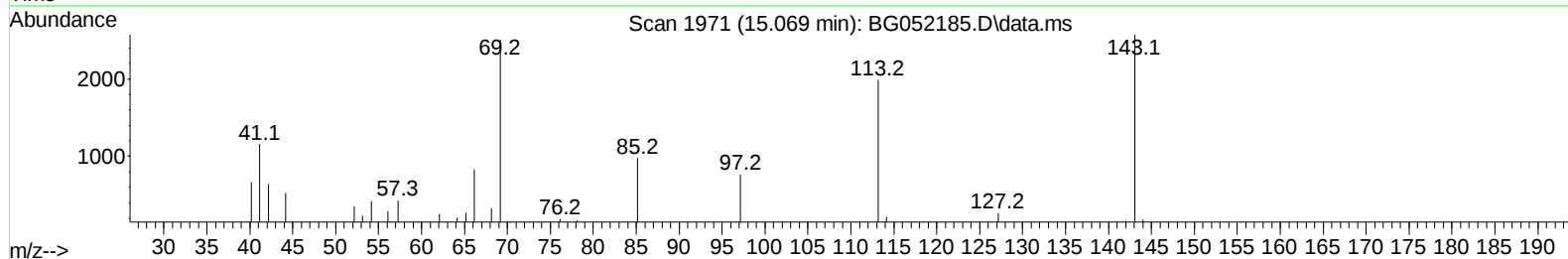
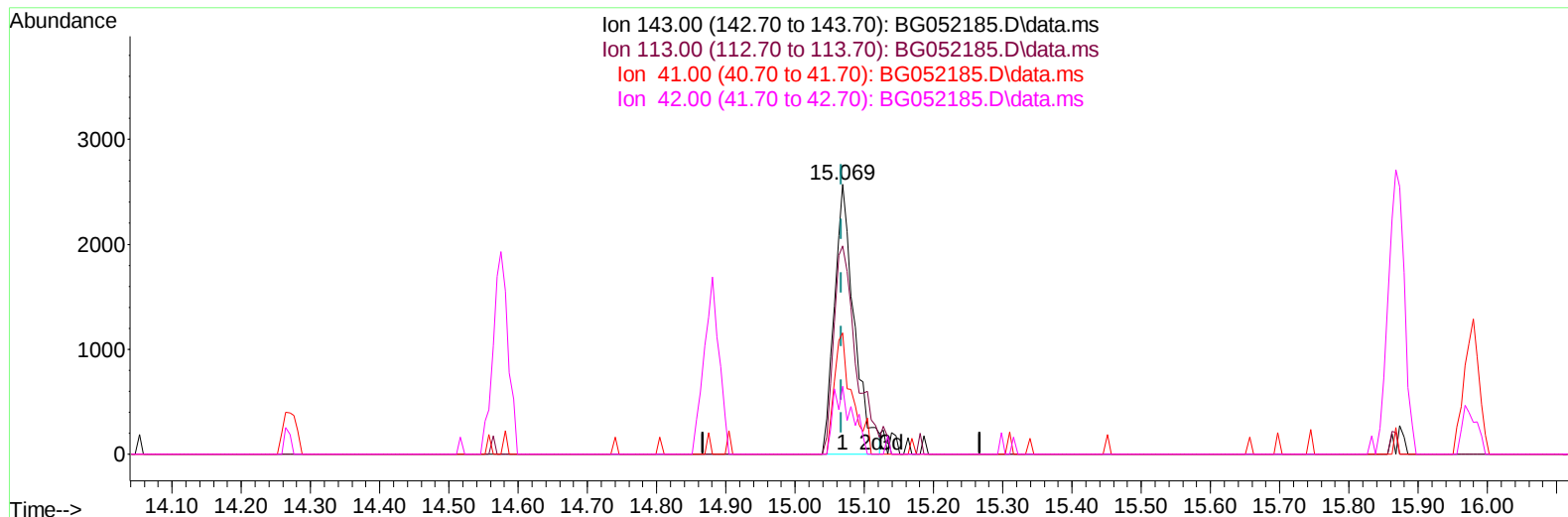
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TIC: BG052185.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.069min (+ 0.002) 5.15 ng/ul m

response 5112

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	77.20
41.00	44.40	45.02
42.00	29.70	25.10

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052185.D
 Acq On : 27 Jan 2022 14:50
 Operator : CG/JU
 Sample : N1279-05
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0VD3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/28/2022
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Quant Time: Jan 27 23:59:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 26 23:58:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.237	152	23100	20.000	ng/ul	0.00
20) Naphthalene-d8	11.074	136	93326	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.881	164	71850	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.630	188	177588	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.948	240	189573	20.000	ng/ul	# 0.00
88) Perylene-d12	25.426	264	187191	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.567	96	2433	3.764	ng/uL	0.00
4) Pyridine-d5	3.996	84	12457	6.199	ng/ul	0.00
7) Phenol-d5	7.379	99	12107	5.355	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.544	67	39749	27.695	ng/ul	0.00
11) 2-Chlorophenol-d4	7.767	132	35190	23.509	ng/ul	0.00
15) 4-Methylphenol-d8	8.942	113	23528	13.387	ng/ul	0.00
21) Nitrobenzene-d5	9.406	128	23742	32.195	ng/ul	0.00
24) 2-Nitrophenol-d4	10.140	143	25218	30.920	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.687	165	44651	28.009	ng/ul	0.00
31) 4-Chloroaniline-d4	11.198	131	54887	24.723	ng/ul	0.00
46) Dimethylphthalate-d6	14.270	166	176643	29.743	ng/ul	0.00
49) Acenaphthylene-d8	14.576	160	219508	30.598	ng/ul	0.00
54) 4-Nitrophenol-d4	15.069	143	5112m	5.148	ng/ul	0.00
60) Fluorene-d10	15.868	176	170527	32.933	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.980	200	25340	22.619	ng/ul	0.00
73) Anthracene-d10	17.730	188	320829	38.083	ng/ul	0.00
81) Pyrene-d10	20.009	212	395920	36.289	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.185	264	377800	38.338	ng/ul	0.00

Target Compounds Qvalue

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

