

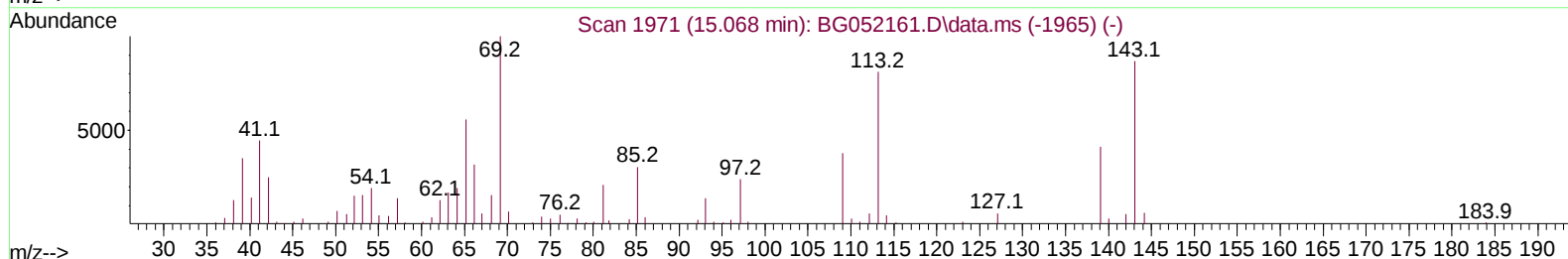
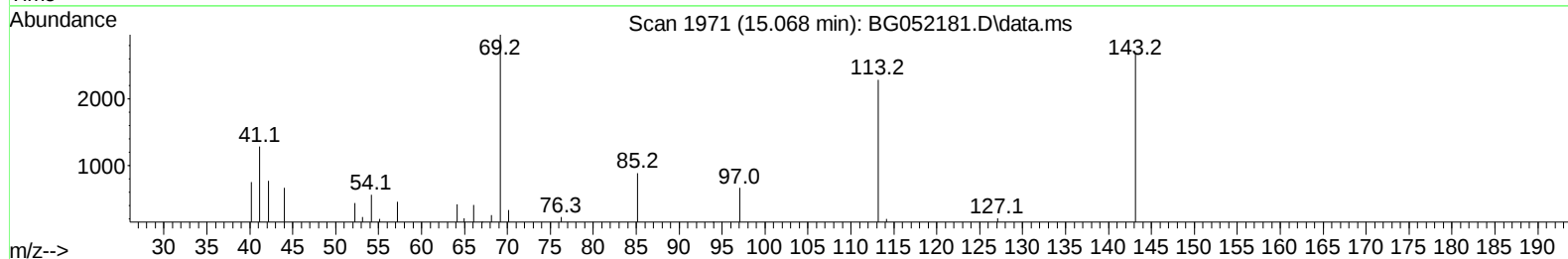
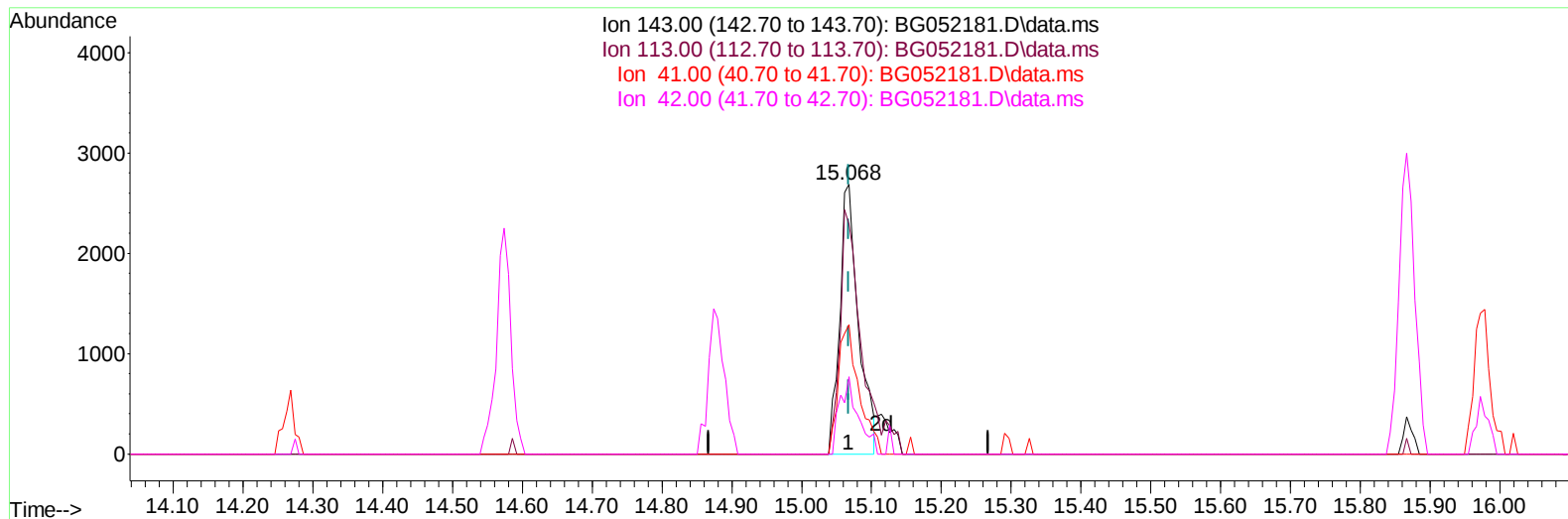
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
Data File : BG052181.D
Acq On : 27 Jan 2022 12:19
Operator : CG/JU
Sample : N1279-01
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0VD8

Manual IntegrationsAPPROVED

Quant Time: Jan 27 23:58:12 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: BG052181.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.068min (+ 0.000) 5.04 ng/u1

response 4981

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	85.11
41.00	44.40	47.92
42.00	29.70	28.67

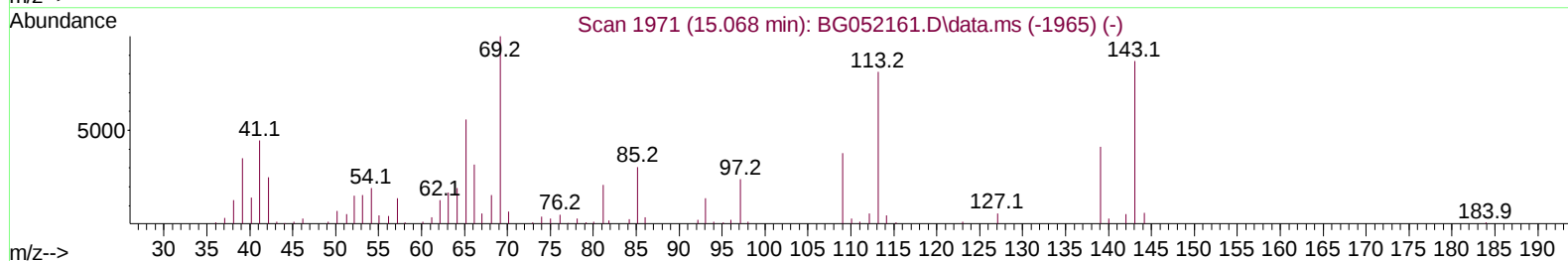
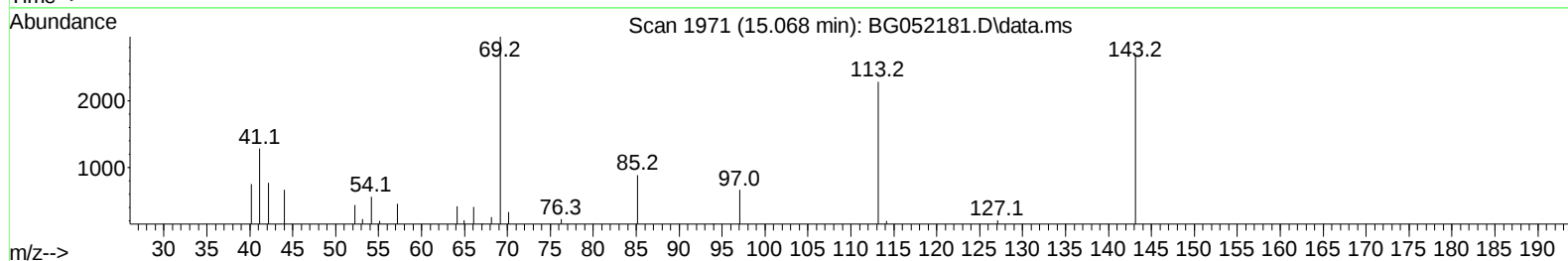
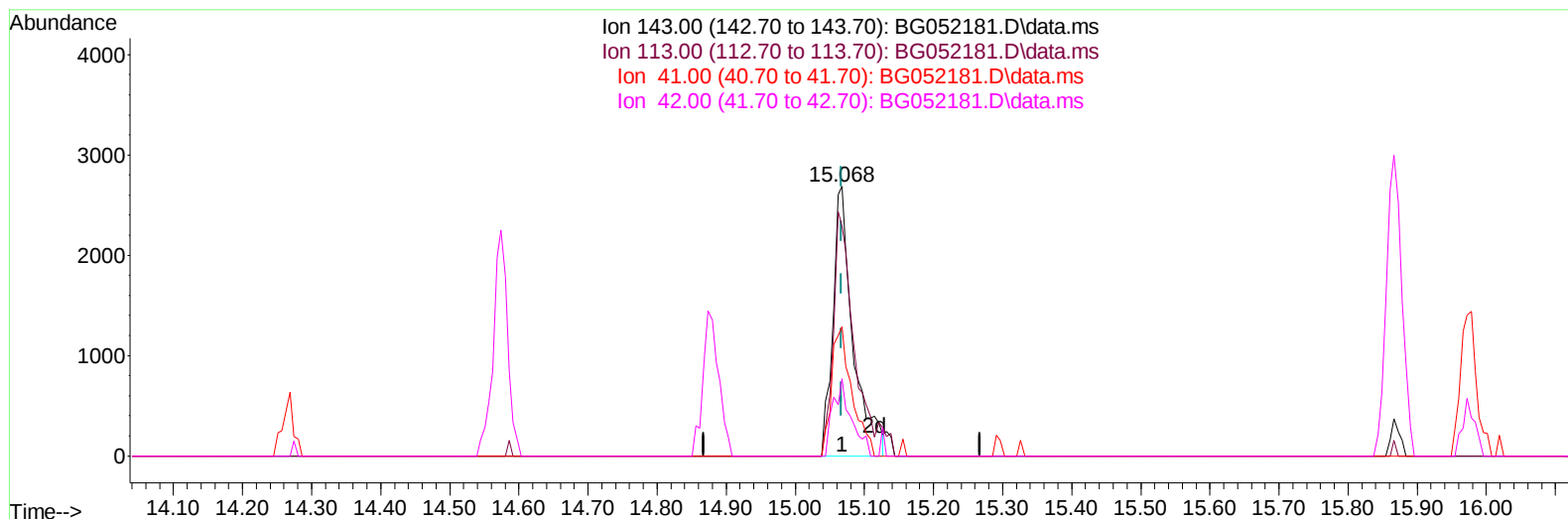
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TIC: BG052181.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.068min (+ 0.000) 5.51 ng/ul m

response 5442

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	85.11
41.00	44.40	47.92
42.00	29.70	28.67

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.236	152	22608	20.000	ng/ul	0.00
20) Naphthalene-d8	11.073	136	93974	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.880	164	71460	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.629	188	178917	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.946	240	185568	20.000	ng/ul	# 0.00
88) Perylene-d12	25.418	264	189019	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.565	96	2661	4.206	ng/uL	0.00
4) Pyridine-d5	4.006	84	2120	1.078	ng/ul	0.01
7) Phenol-d5	7.378	99	13567	6.131	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	42589	30.320	ng/ul	0.00
11) 2-Chlorophenol-d4	7.766	132	37280	25.447	ng/ul	0.00
15) 4-Methylphenol-d8	8.941	113	25027	14.550	ng/ul	0.00
21) Nitrobenzene-d5	9.405	128	25716	34.632	ng/ul	0.00
24) 2-Nitrophenol-d4	10.139	143	27539	33.533	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.691	165	49411	30.781	ng/ul	0.00
31) 4-Chloroaniline-d4	11.202	131	33095	14.804	ng/ul	0.00
46) Dimethylphthalate-d6	14.269	166	199056	33.700	ng/ul	0.00
49) Acenaphthylene-d8	14.574	160	240779	33.746	ng/ul	0.00
54) 4-Nitrophenol-d4	15.068	143	5442m	5.511	ng/ul	0.00
60) Fluorene-d10	15.866	176	177954	34.555	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.978	200	26245	23.253	ng/ul	0.00
73) Anthracene-d10	17.729	188	317543	37.413	ng/ul	0.00
81) Pyrene-d10	20.008	212	399217	37.381	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.183	264	374232	37.608	ng/ul	0.00

Target Compounds Qvalue

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

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