

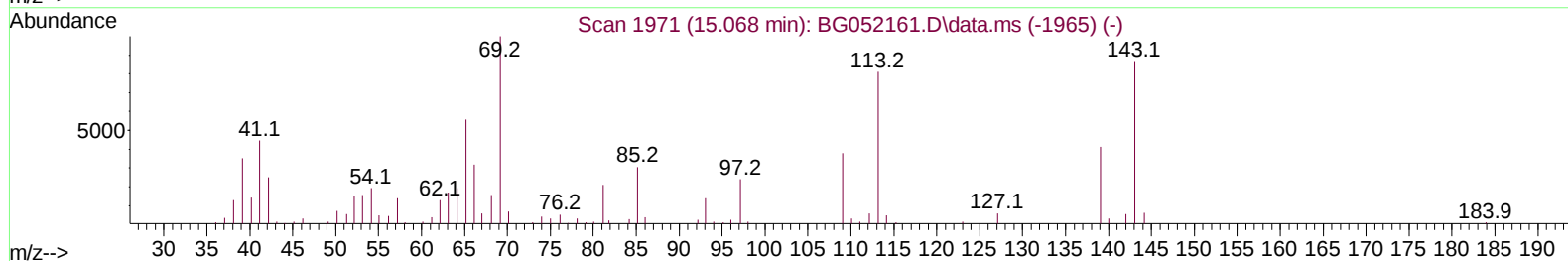
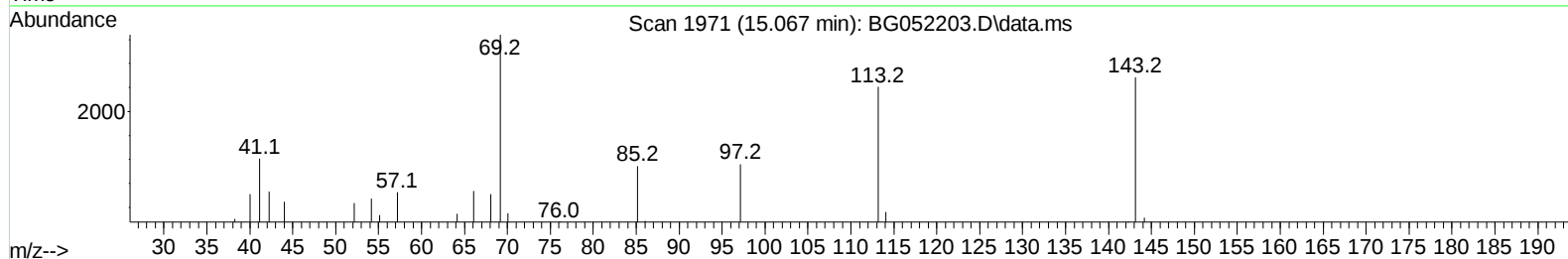
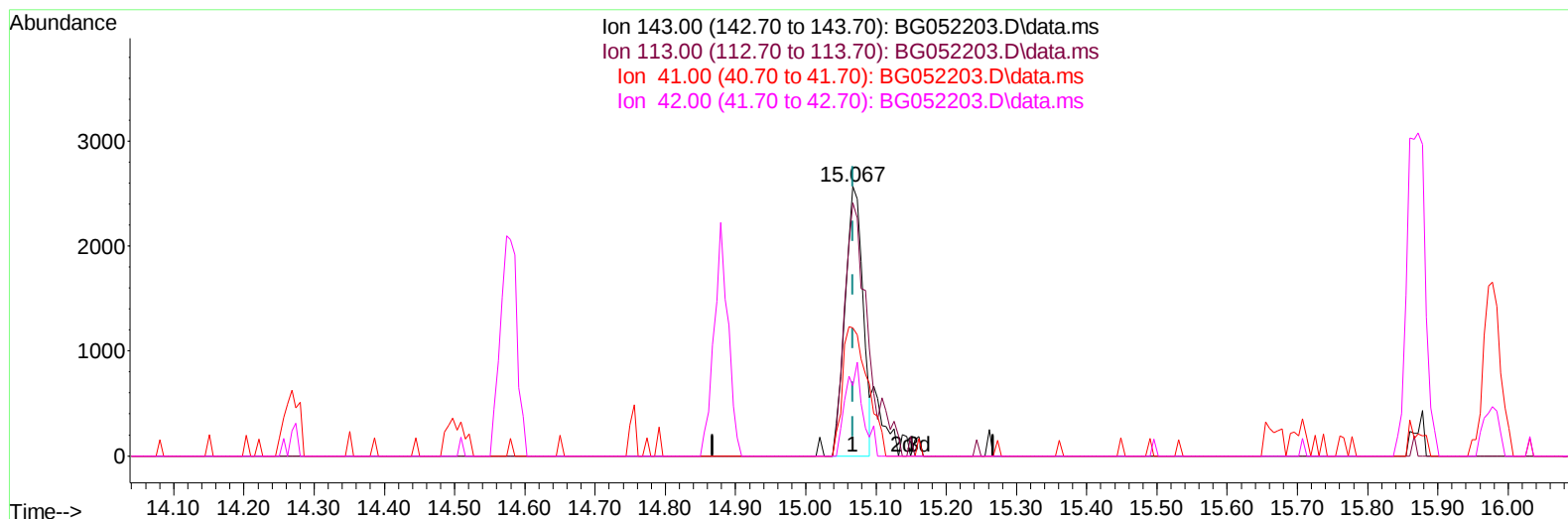
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
Data File : BG052203.D
Acq On : 28 Jan 2022 2:46
Operator : CG/JU
Sample : N1302-03
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0F33

Manual IntegrationsAPPROVED

Quant Time: Jan 28 04:08:47 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: BG052203.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.067min (0.000) 4.17 ng/u1

response 4559

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	93.92
41.00	44.40	47.45
42.00	29.70	25.98

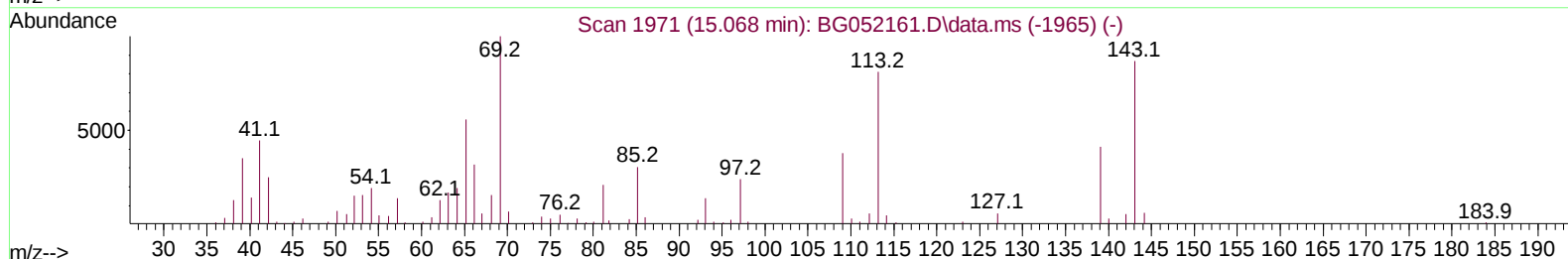
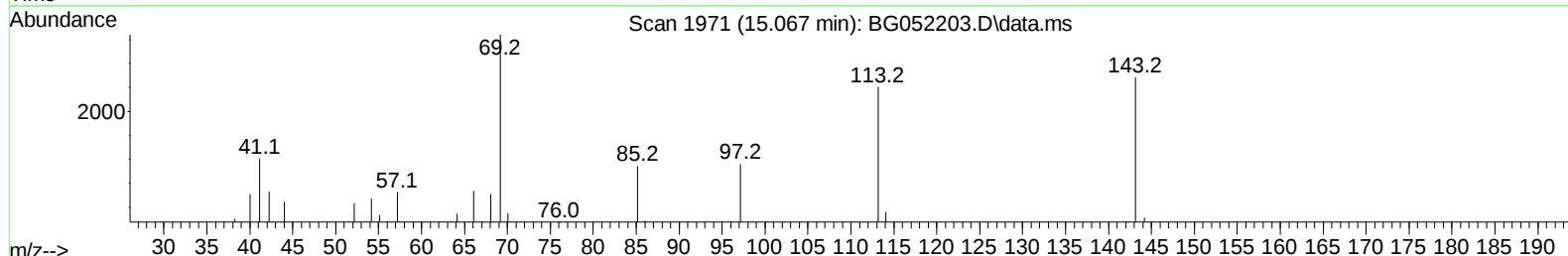
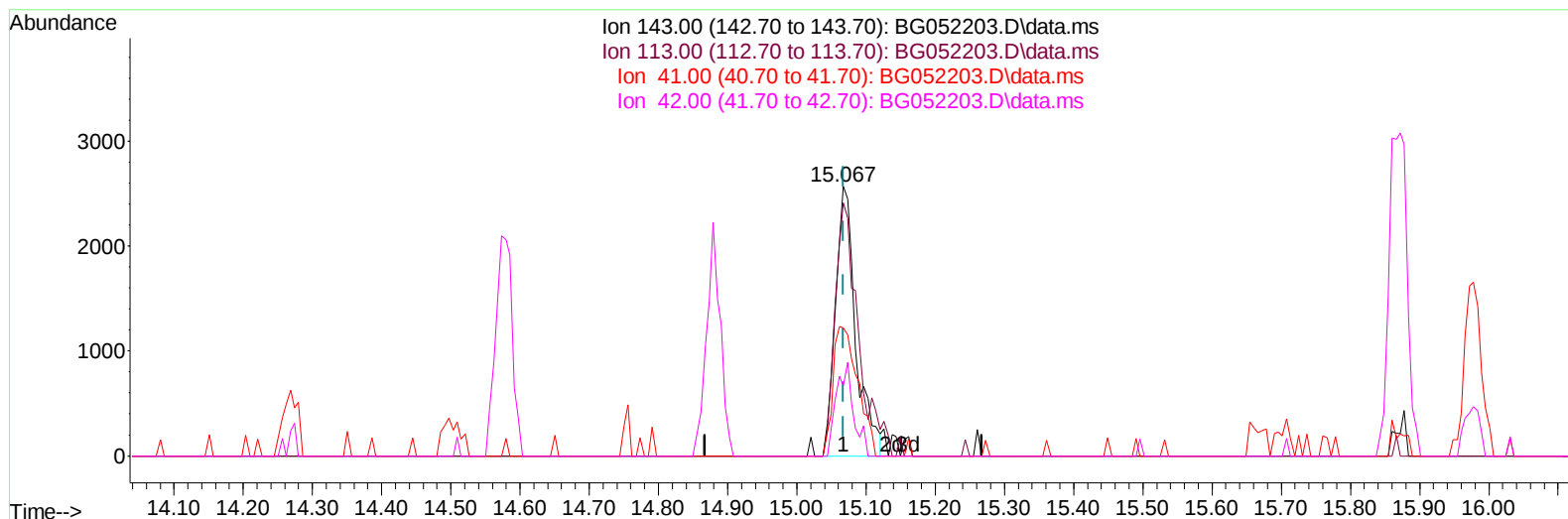
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TIC: BG052203.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.067min (0.000) 4.81 ng/ul m

response 5258

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	93.92
41.00	44.40	47.45
42.00	29.70	25.98

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.235	152	24810	20.000	ng/ul	0.00
20) Naphthalene-d8	11.078	136	105706	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.879	164	79170	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.628	188	184639	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.946	240	193042	20.000	ng/ul	# 0.00
88) Perylene-d12	25.430	264	198578	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.571	96	2852	4.108	ng/uL	0.00
4) Pyridine-d5	4.000	84	17046	7.898	ng/ul	0.00
7) Phenol-d5	7.378	99	16113	6.636	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	51907	33.674	ng/ul	0.00
11) 2-Chlorophenol-d4	7.765	132	43350	26.964	ng/ul	0.00
15) 4-Methylphenol-d8	8.940	113	29442	15.597	ng/ul	0.00
21) Nitrobenzene-d5	9.410	128	31437	37.637	ng/ul	0.00
24) 2-Nitrophenol-d4	10.139	143	33472	36.234	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.691	165	58798	32.563	ng/ul	0.00
31) 4-Chloroaniline-d4	11.202	131	79009	31.420	ng/ul	0.00
46) Dimethylphthalate-d6	14.268	166	233780	35.724	ng/ul	0.00
49) Acenaphthylene-d8	14.580	160	289392	36.610	ng/ul	0.00
54) 4-Nitrophenol-d4	15.067	143	5258m	4.806	ng/ul	0.00
60) Fluorene-d10	15.872	176	209713	36.756	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.978	200	28517	24.483	ng/ul	0.00
73) Anthracene-d10	17.728	188	350221	39.985	ng/ul	0.00
81) Pyrene-d10	20.008	212	422896	38.065	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.189	264	419316	40.110	ng/ul	0.00

Target Compounds Qvalue

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

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