

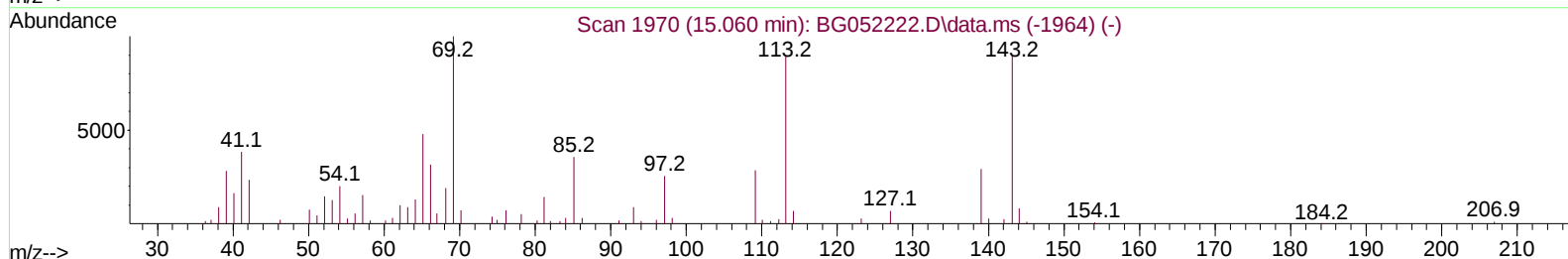
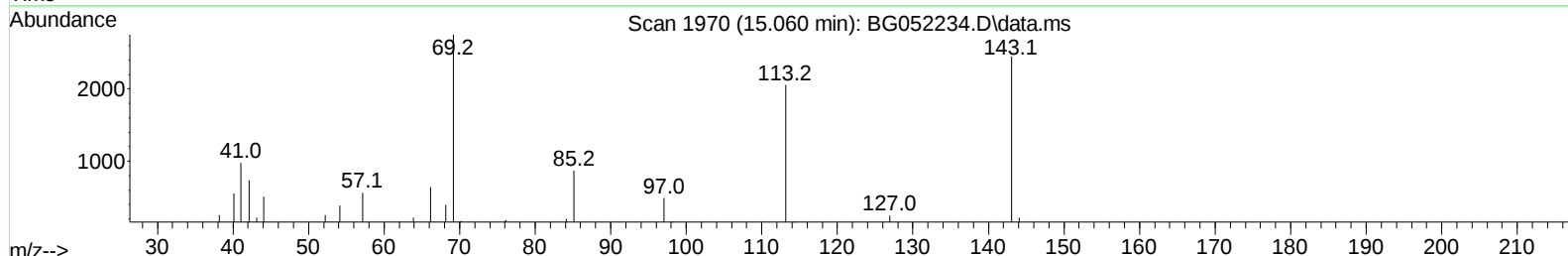
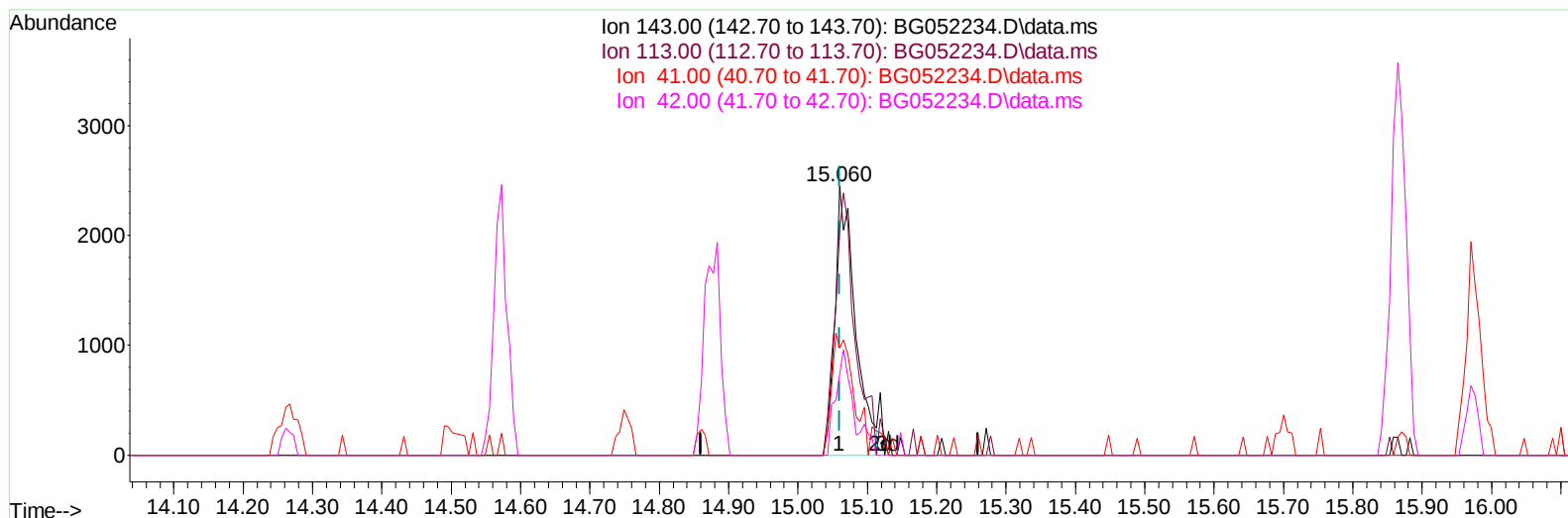
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020122\
 Data File : BG052234.D
 Acq On : 1 Feb 2022 17:49
 Operator : CG/JU
 Sample : N1328-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COHM7

Manual IntegrationsAPPROVED

Quant Time: Feb 02 02:46:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG013122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jan 31 18:07:10 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/03/2022
 Supervised By :mohammad ahmed 02/04/2022



TIC: BG052234.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.060min (-0.000) 4.74 ng/u1

response 4975

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	83.87
41.00	44.40	39.85
42.00	29.70	30.09

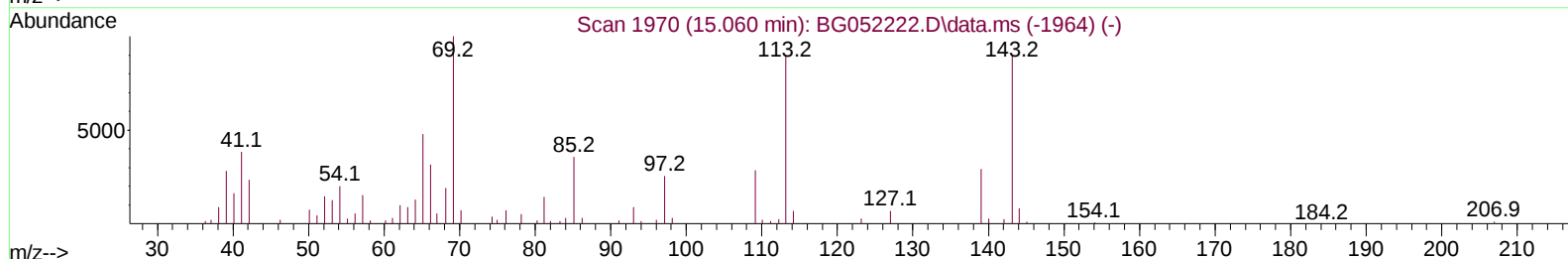
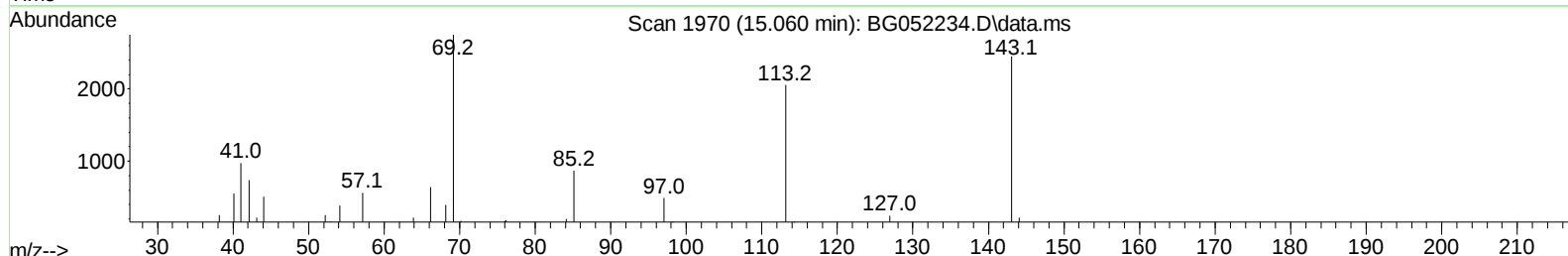
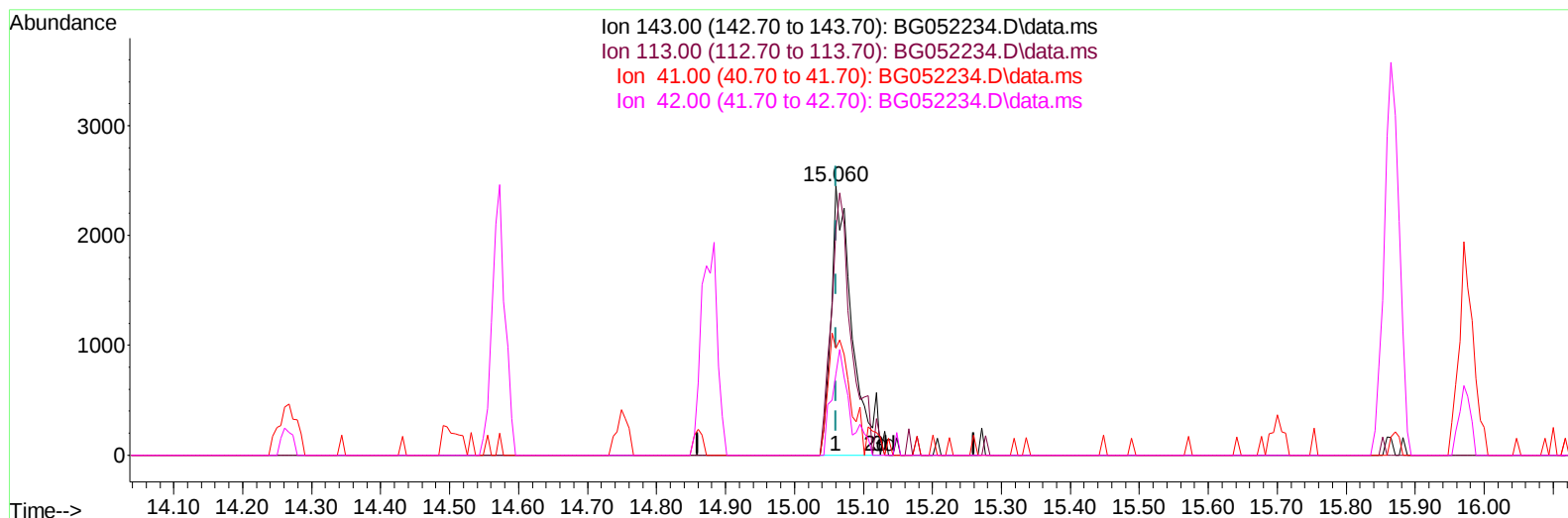
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TIC: BG052234.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.060min (-0.000) 4.93 ng/u1 m

response 5177

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	83.87
41.00	44.40	39.85
42.00	29.70	30.09

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.234	152	29838	20.000	ng/ul	0.00
20) Naphthalene-d8	11.071	136	128075	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.878	164	88142	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.627	188	212607	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.939	240	212064	20.000	ng/ul	# 0.00
88) Perylene-d12	25.417	264	213825	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.564	96	2820	3.813	ng/uL	0.00
4) Pyridine-d5	3.993	84	14156	6.154	ng/ul	0.00
7) Phenol-d5	7.376	99	15257	5.621	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.535	67	52362	31.757	ng/ul	0.00
11) 2-Chlorophenol-d4	7.764	132	44295	23.421	ng/ul	0.00
15) 4-Methylphenol-d8	8.933	113	28627	13.899	ng/ul	0.00
21) Nitrobenzene-d5	9.403	128	30963	29.691	ng/ul	0.00
24) 2-Nitrophenol-d4	10.137	143	31720	27.836	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.689	165	53399	24.037	ng/ul	0.00
31) 4-Chloroaniline-d4	11.201	131	66794	23.704	ng/ul	0.00
46) Dimethylphthalate-d6	14.267	166	223041	32.193	ng/ul	0.00
49) Acenaphthylene-d8	14.572	160	267388	30.963	ng/ul	0.00
54) 4-Nitrophenol-d4	15.060	143	5177m	4.934	ng/ul	0.00
60) Fluorene-d10	15.865	176	198460	32.696	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.971	200	32954	25.106	ng/ul	-0.01
73) Anthracene-d10	17.721	188	340042	33.926	ng/ul	0.00
81) Pyrene-d10	20.006	212	413216	32.512	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.176	264	382180	33.643	ng/ul	0.00

Target Compounds Qvalue

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

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