

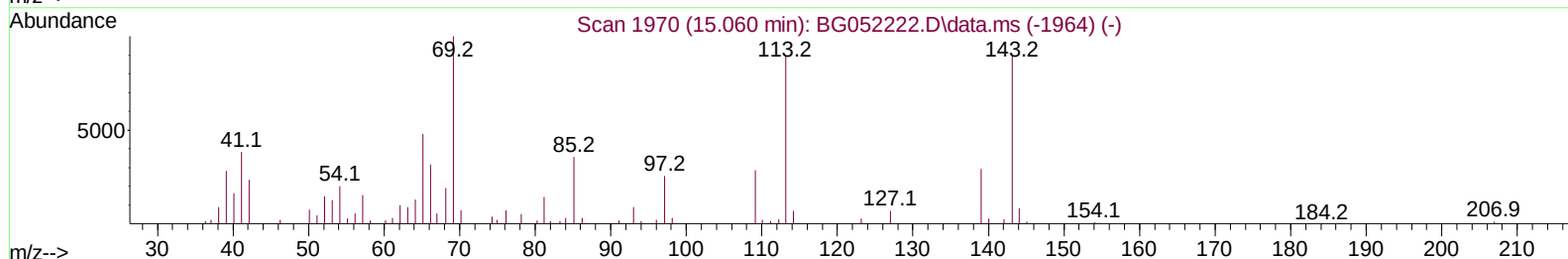
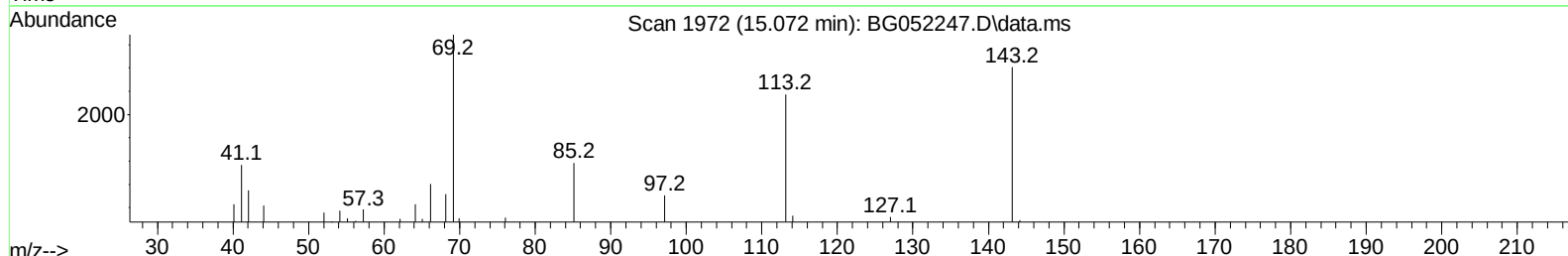
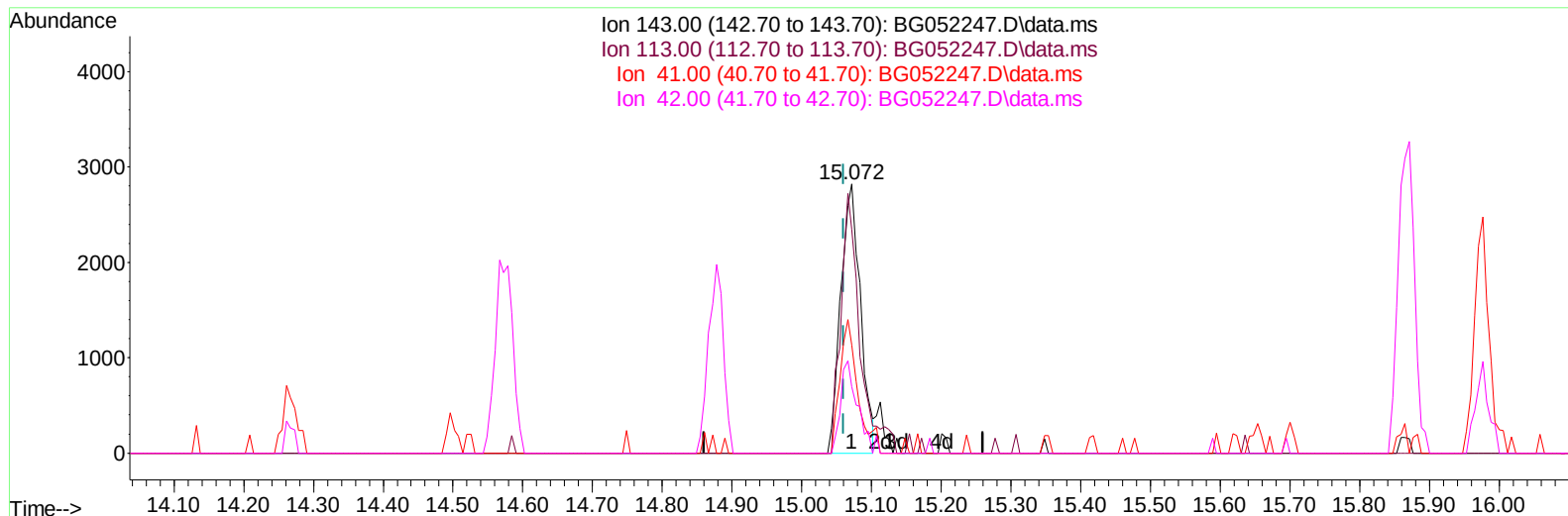
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020122\
Data File : BG052247.D
Acq On : 2 Feb 2022 2:37
Operator : CG/JU
Sample : N1307-11
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0Q34

Manual IntegrationsAPPROVED

Quant Time: Feb 02 03:52:58 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: BG052247.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.072min (+ 0.012) 5.13 ng/ul

response 5495

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	83.23
41.00	44.40	40.28
42.00	29.70	24.61

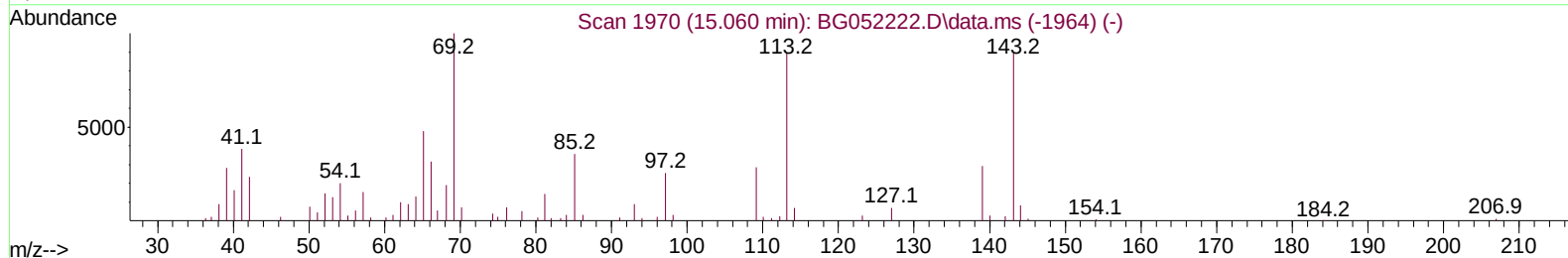
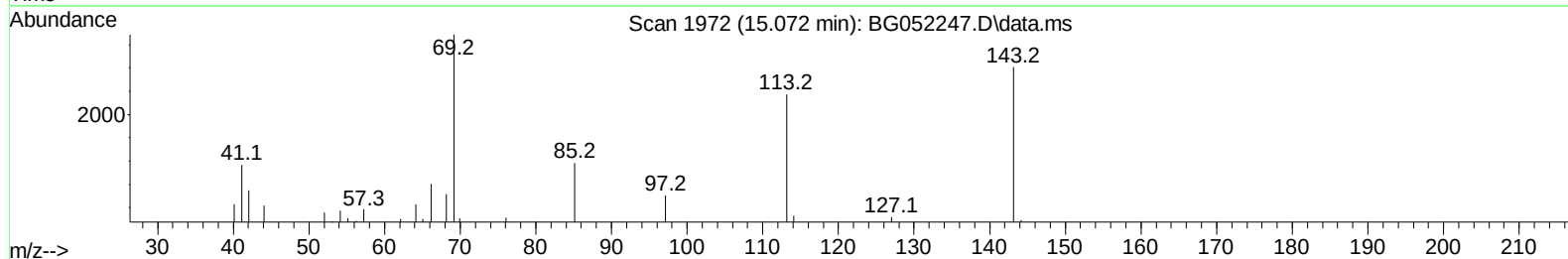
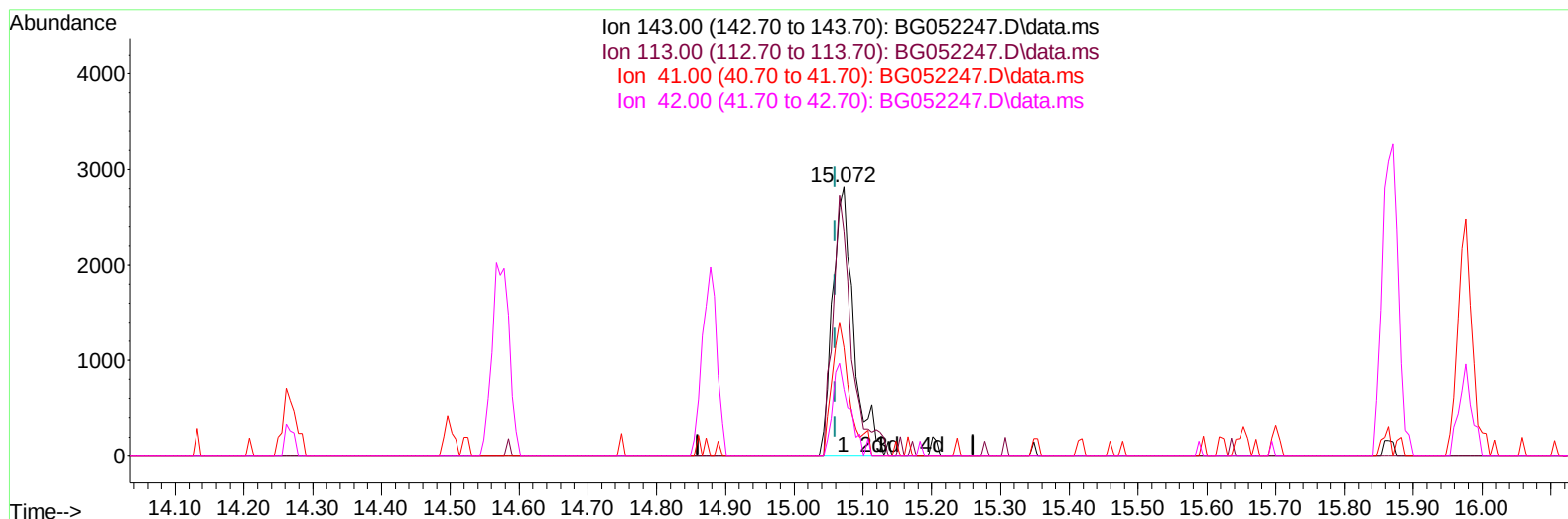
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TIC: BG052247.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.072min (+ 0.012) 5.55 ng/ul m

response 5946

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	83.23
41.00	44.40	40.28
42.00	29.70	24.61

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.234	152	30466	20.000	ng/ul	0.00
20) Naphthalene-d8	11.077	136	126390	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.878	164	90015	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.627	188	214945	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.939	240	210373	20.000	ng/ul	# 0.00
88) Perylene-d12	25.417	264	213009	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.564	96	2966	3.928	ng/uL	0.00
4) Pyridine-d5	3.993	84	19450	8.281	ng/ul	0.00
7) Phenol-d5	7.376	99	16216	5.852	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	53681	31.885	ng/ul	0.00
11) 2-Chlorophenol-d4	7.764	132	44788	23.193	ng/ul	0.00
15) 4-Methylphenol-d8	8.939	113	31395	14.929	ng/ul	0.00
21) Nitrobenzene-d5	9.403	128	32063	31.156	ng/ul	0.00
24) 2-Nitrophenol-d4	10.137	143	34509	30.687	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.690	165	58399	26.639	ng/ul	0.00
31) 4-Chloroaniline-d4	11.201	131	71062	25.554	ng/ul	0.00
46) Dimethylphthalate-d6	14.267	166	235445	33.276	ng/ul	0.00
49) Acenaphthylene-d8	14.573	160	288449	32.707	ng/ul	0.00
54) 4-Nitrophenol-d4	15.072	143	5946m	5.549	ng/ul	0.01
60) Fluorene-d10	15.865	176	208936	33.706	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.977	200	36909	27.814	ng/ul	0.00
73) Anthracene-d10	17.727	188	350509	34.590	ng/ul	0.00
81) Pyrene-d10	20.006	212	417710	33.130	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.176	264	389384	34.409	ng/ul	0.00

Target Compounds Qvalue

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

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