

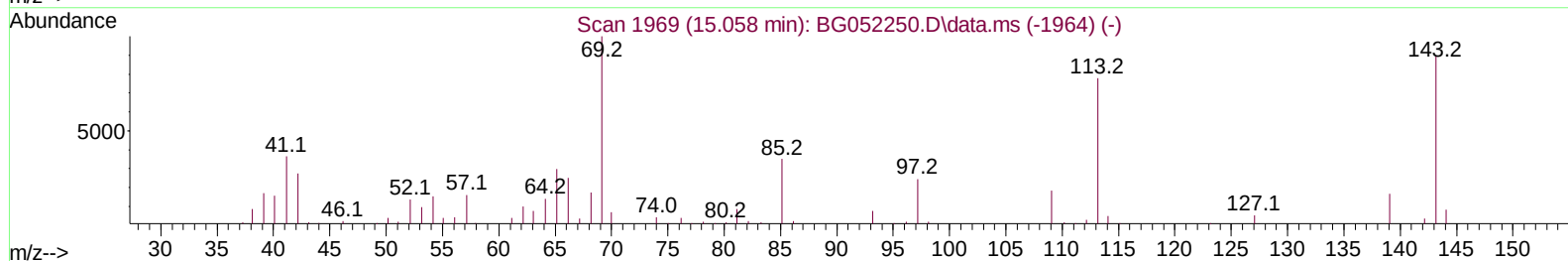
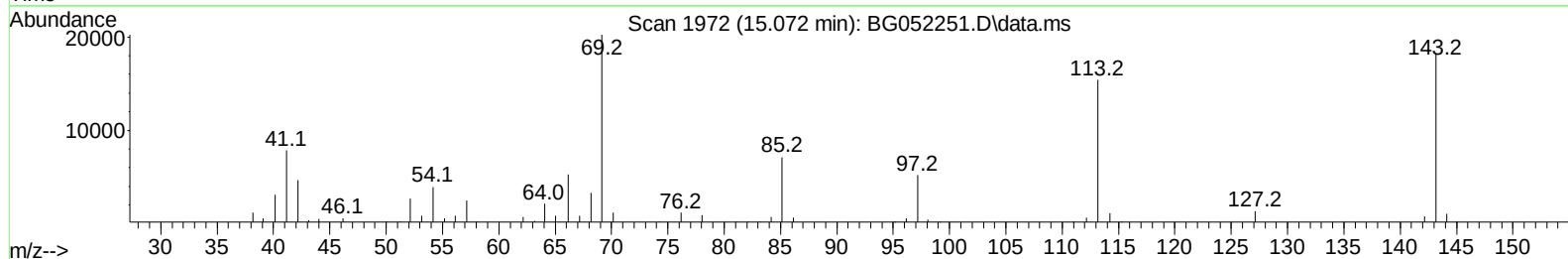
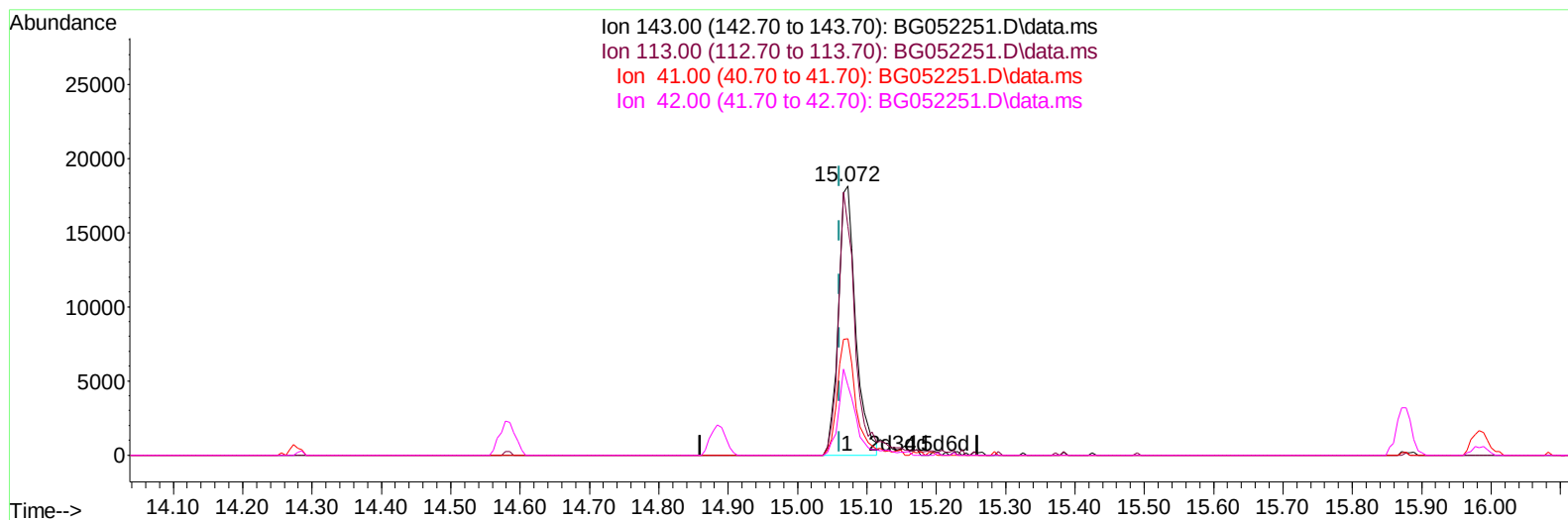
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020122\
 Data File : BG052251.D
 Acq On : 2 Feb 2022 8:49
 Operator : CG/JU
 Sample : PB142387BL
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK387

Manual IntegrationsAPPROVED

Quant Time: Feb 03 04:02:56 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: BG052251.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.072min (+ 0.012) 28.06 ng/ul

response 31428

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	84.95
41.00	44.40	43.37
42.00	29.70	25.80

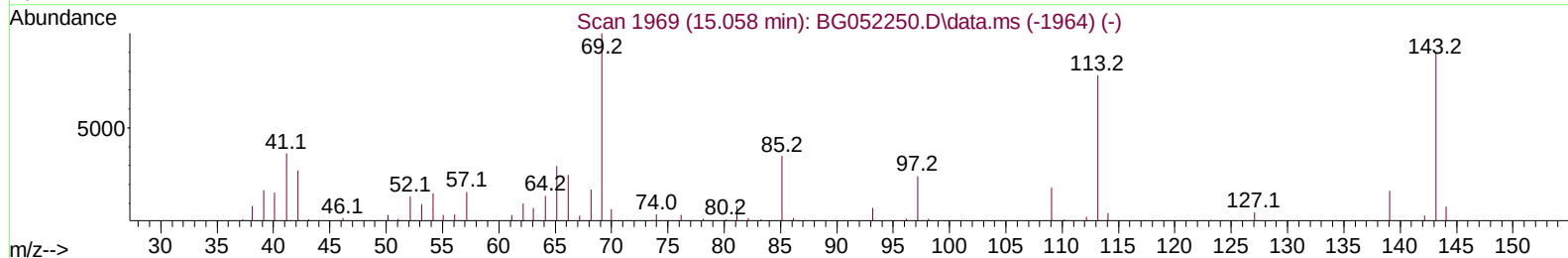
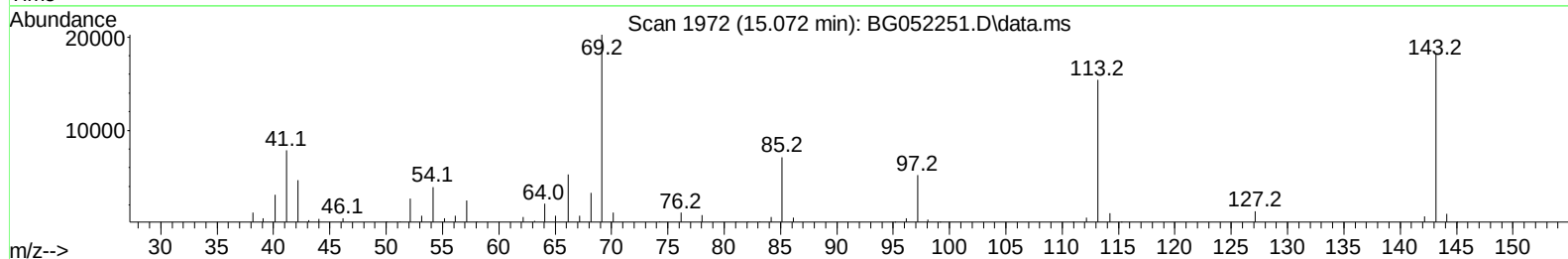
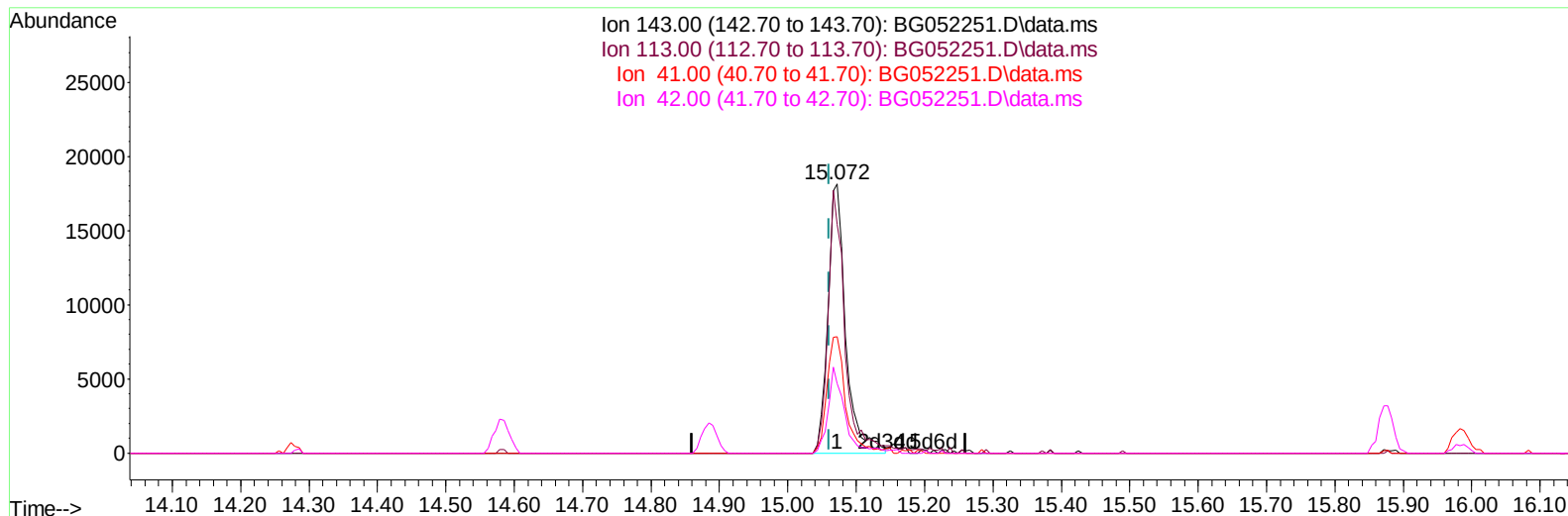
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TIC: BG052251.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.072min (+ 0.012) 29.11 ng/ul m

response 32605

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	84.95
41.00	44.40	43.37
42.00	29.70	25.80

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.252	152	30237	20.000	ng/ul	0.02
20) Naphthalene-d8	11.090	136	129907	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.890	164	94080	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.634	188	223688	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.946	240	223492	20.000	ng/ul	# 0.00
88) Perylene-d12	25.429	264	224023	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.570	96	4524	6.036	ng/uL	0.00
4) Pyridine-d5	3.999	84	68770	29.500	ng/ul	0.01
7) Phenol-d5	7.395	99	81132	29.499	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.559	67	52446	31.388	ng/ul	0.02
11) 2-Chlorophenol-d4	7.776	132	59297	30.939	ng/ul	0.01
15) 4-Methylphenol-d8	8.951	113	64640	30.971	ng/ul	0.01
21) Nitrobenzene-d5	9.421	128	32977	31.177	ng/ul	0.01
24) 2-Nitrophenol-d4	10.156	143	35830	30.999	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.702	165	64398	28.580	ng/ul	0.01
31) 4-Chloroaniline-d4	11.213	131	84239	29.473	ng/ul	0.01
46) Dimethylphthalate-d6	14.279	166	234497	31.710	ng/ul	0.01
49) Acenaphthylene-d8	14.585	160	276803	30.031	ng/ul	0.00
54) 4-Nitrophenol-d4	15.072	143	32605m	29.112	ng/ul	0.01
60) Fluorene-d10	15.877	176	196498	30.330	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.989	200	36048	26.103	ng/ul	0.00
73) Anthracene-d10	17.734	188	350287	33.217	ng/ul	0.00
81) Pyrene-d10	20.013	212	425256	31.749	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.188	264	400146	33.622	ng/ul	0.00

Target Compounds Qvalue

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

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