

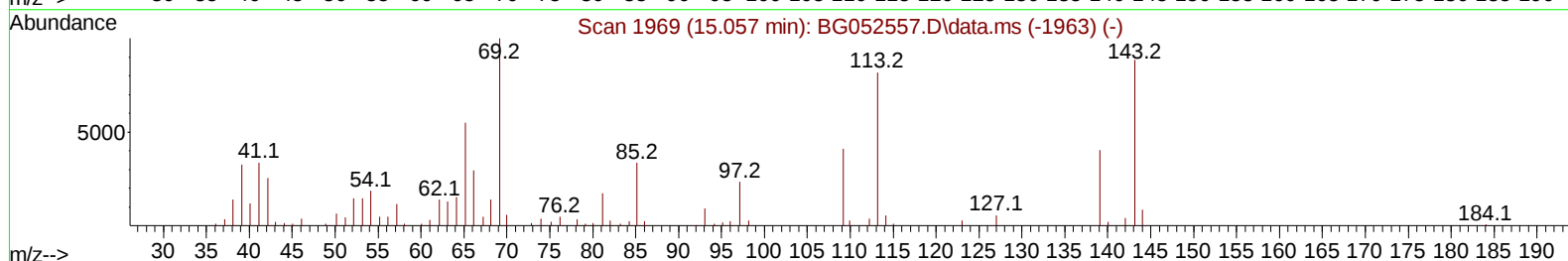
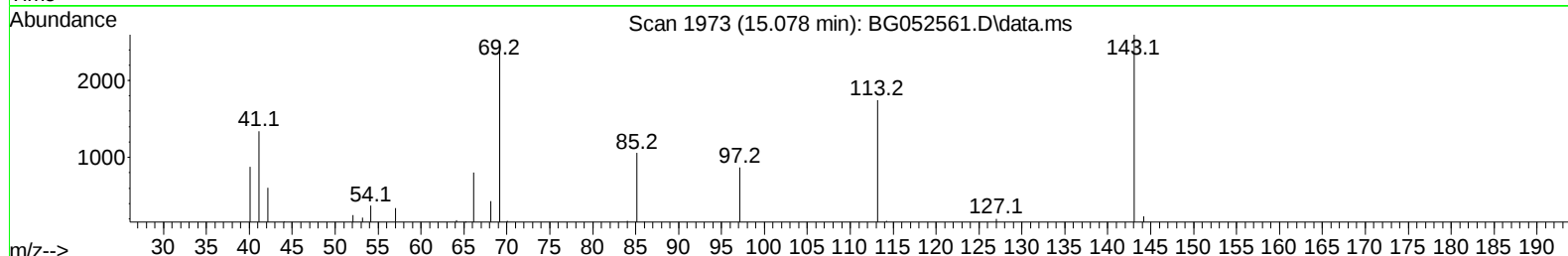
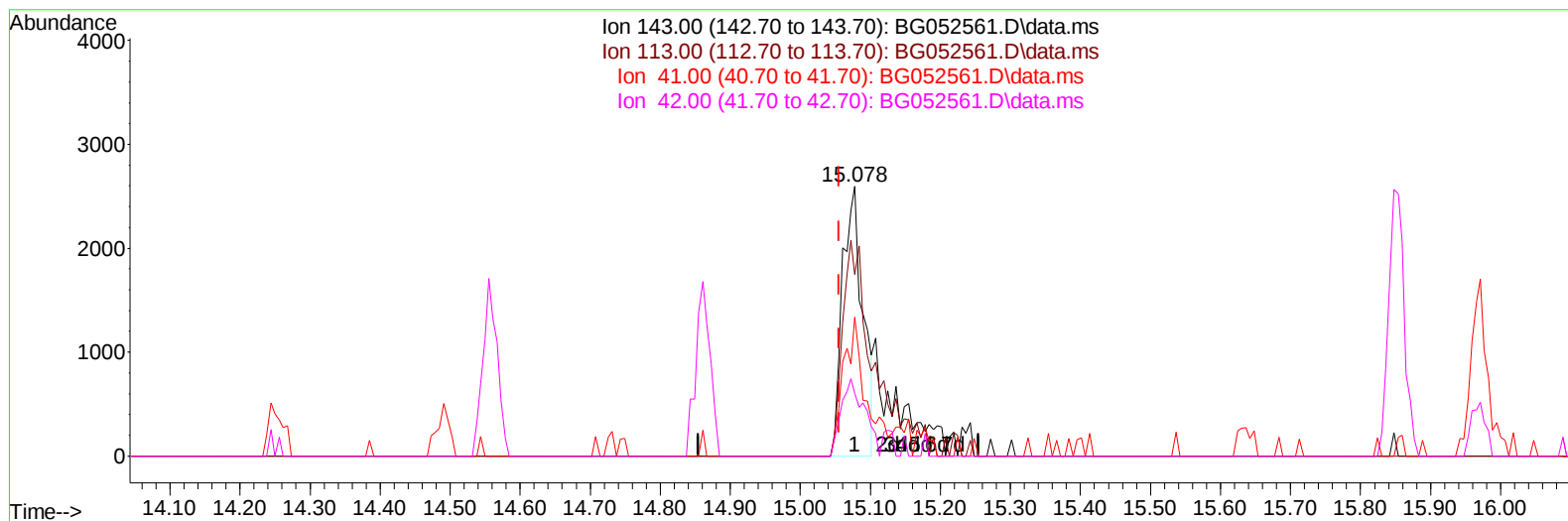
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030322\
 Data File : BG052561.D
 Acq On : 3 Mar 2022 19:18
 Operator : CG/JU
 Sample : N1745-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 H0AA3

Manual IntegrationsAPPROVED

Quant Time: Mar 04 00:02:52 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022222.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 22 17:32:12 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/04/2022
 Supervised By :Yogesh Patel 03/07/2022



TIC: BG052561.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.078min (+ 0.023) 5.41 ng/ul

response 5326

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	67.35
41.00	44.40	51.66
42.00	29.70	23.17#

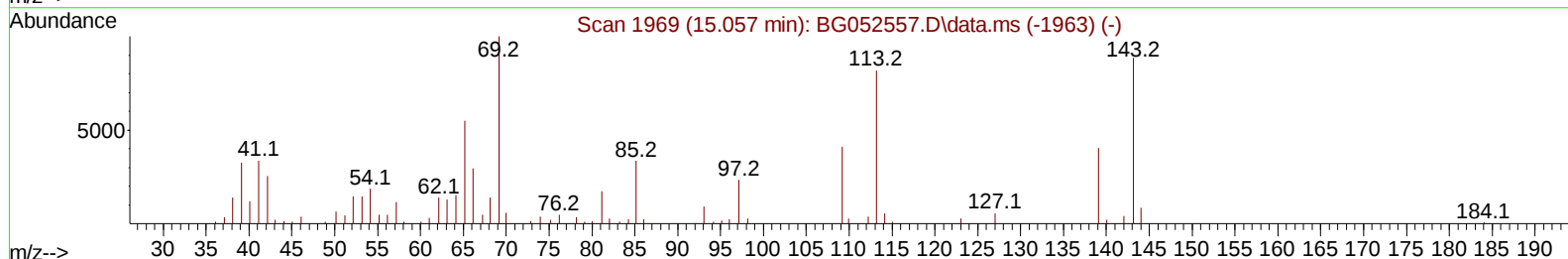
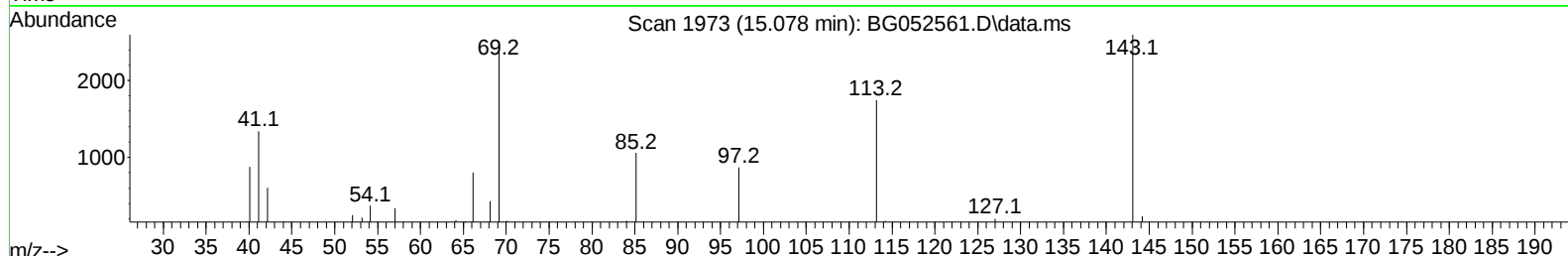
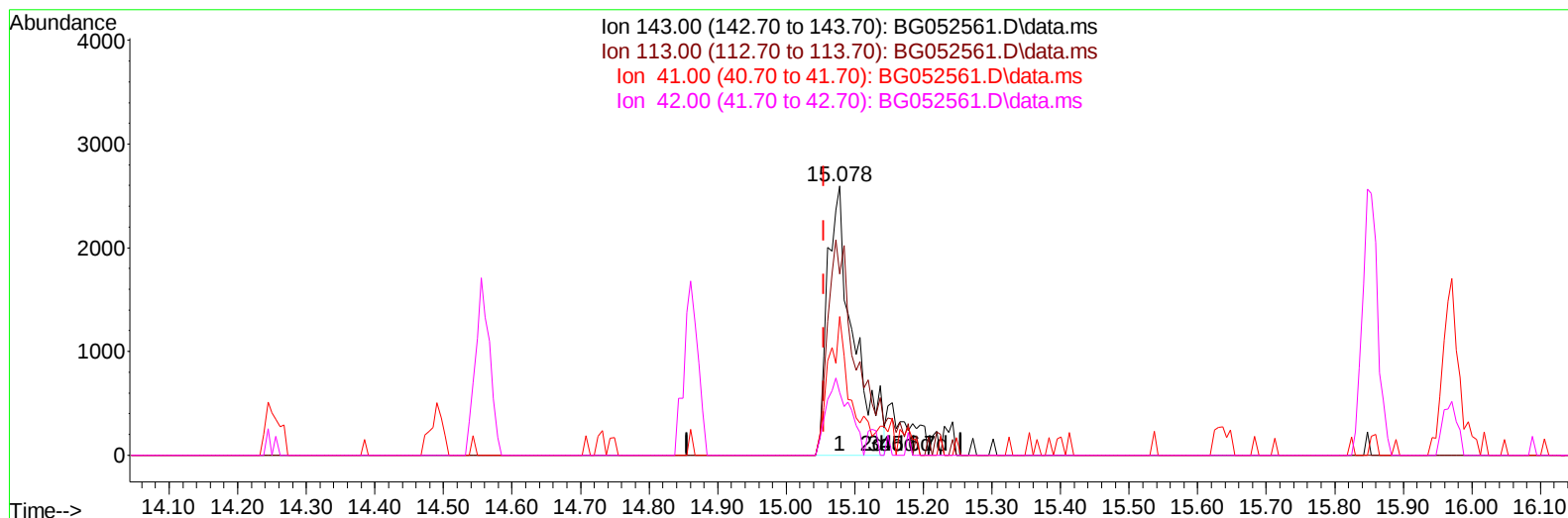
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TIC: BG052561.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.078min (+ 0.023) 6.87 ng/ul m

response 6763

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	67.35
41.00	44.40	51.66
42.00	29.70	23.17#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.223	152	22181	20.000	ng/ul	0.00
20) Naphthalene-d8	11.054	136	99331	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.861	164	76169	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.610	188	193820	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.928	240	200560	20.000	ng/ul	# 0.00
88) Perylene-d12	25.388	264	198367	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.541	96	2548	4.506	ng/uL	0.00
4) Pyridine-d5	3.987	84	10038	5.885	ng/ul	0.01
7) Phenol-d5	7.365	99	16215	7.709	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.524	67	39078	31.792	ng/ul	0.00
11) 2-Chlorophenol-d4	7.747	132	36175	24.487	ng/ul	0.00
15) 4-Methylphenol-d8	8.928	113	28093	17.167	ng/ul	0.00
21) Nitrobenzene-d5	9.392	128	24466	28.893	ng/ul	0.00
24) 2-Nitrophenol-d4	10.120	143	26950	29.056	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.672	165	44956	25.991	ng/ul	0.00
31) 4-Chloroaniline-d4	11.183	131	55713	23.243	ng/ul	0.00
46) Dimethylphthalate-d6	14.250	166	201309	32.511	ng/ul	0.00
49) Acenaphthylene-d8	14.555	160	239889	31.383	ng/ul	0.00
54) 4-Nitrophenol-d4	15.078	143	6763m	6.871	ng/ul	0.02
60) Fluorene-d10	15.854	176	178227	32.765	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.971	200	33909	27.592	ng/ul	0.00
73) Anthracene-d10	17.710	188	342766	36.607	ng/ul	0.00
81) Pyrene-d10	19.989	212	431900	36.802	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.153	264	416691	39.155	ng/ul	0.00

Target Compounds Qvalue

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

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