

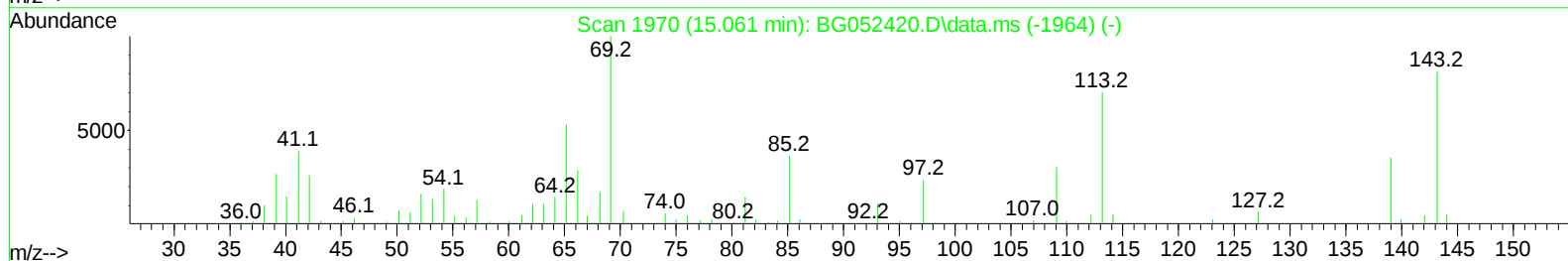
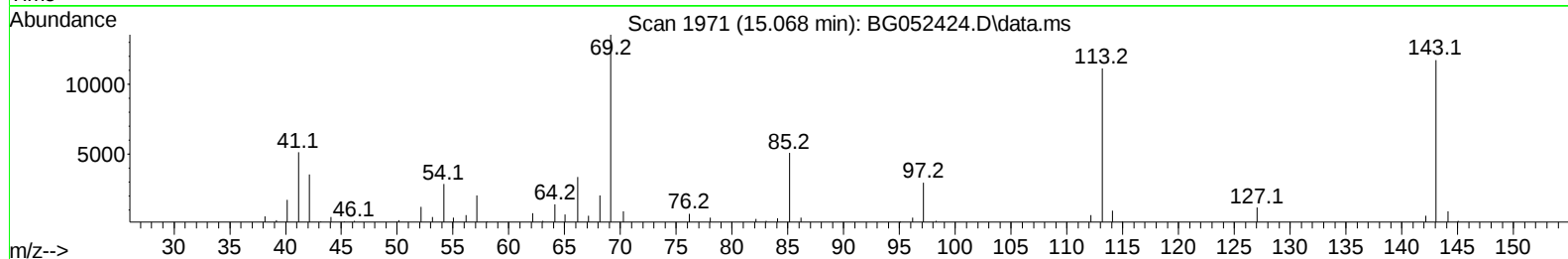
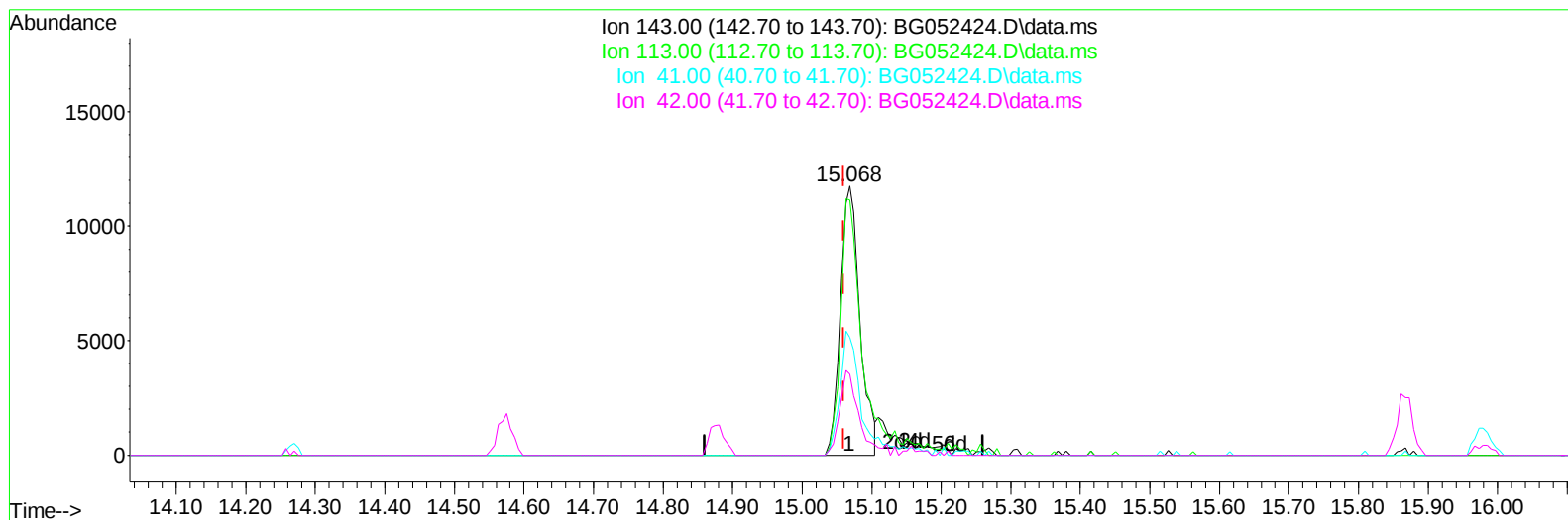
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021822\
Data File : BG052424.D
Acq On : 18 Feb 2022 16:15
Operator : CG/JU
Sample : PB142783BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK783

Manual IntegrationsAPPROVED

Quant Time: Feb 18 22:35:25 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 02/21/2022
Supervised By :Yogesh Patel 02/22/2022



TIC: BG052424.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.068min (+ 0.008) 26.21 ng/u1

response 22955

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	94.85
41.00	44.40	43.66
42.00	29.70	30.11

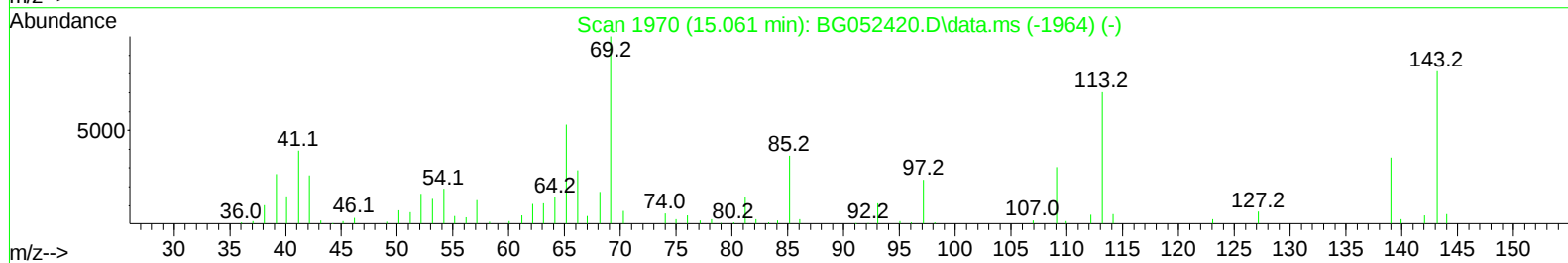
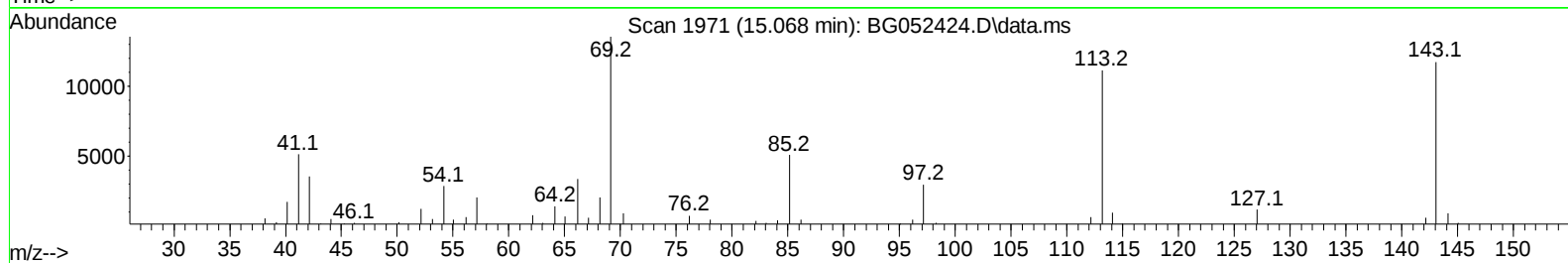
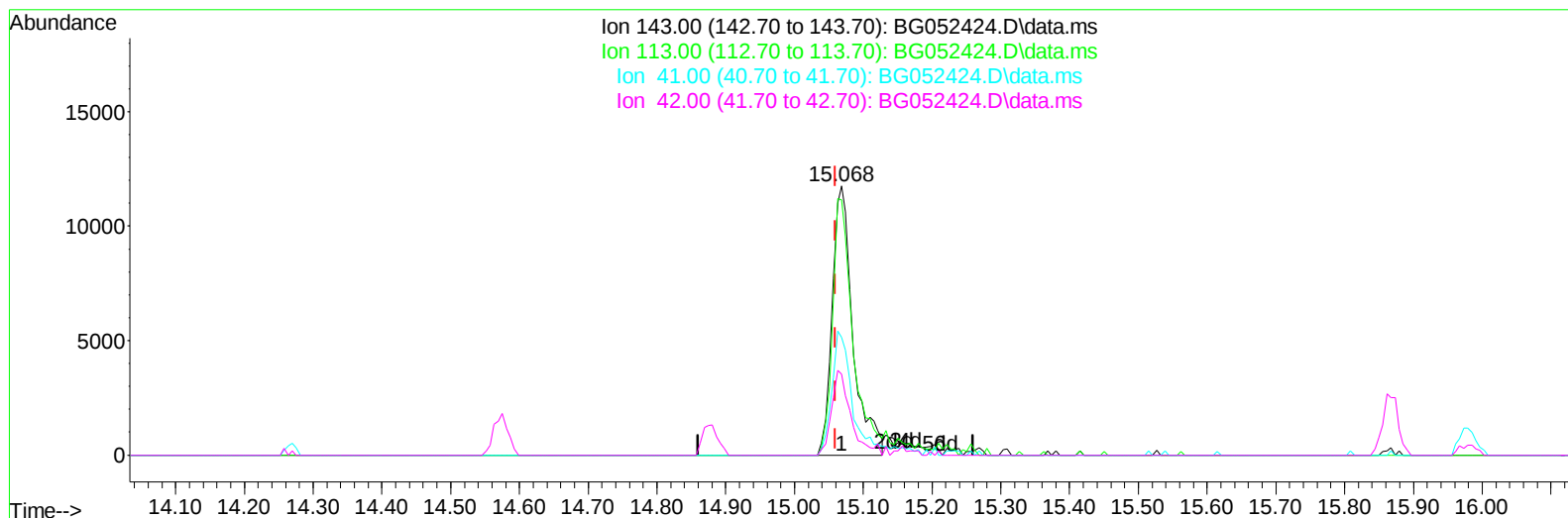
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TIC: BG052424.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.068min (+ 0.008) 28.17 ng/ul m

response 24675

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	94.85
41.00	44.40	43.66
42.00	29.70	30.11

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.237	152	24370	20.000	ng/ul	0.00
20) Naphthalene-d8	11.074	136	100728	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.881	164	73576	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.624	188	182204	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.942	240	204379	20.000	ng/ul	# 0.00
88) Perylene-d12	25.413	264	196278	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.555	96	3838	6.354	ng/uL	0.00
4) Pyridine-d5	3.983	84	59259	31.540	ng/ul	0.00
7) Phenol-d5	7.385	99	68479	30.892	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.543	67	43483	32.289	ng/ul	0.00
11) 2-Chlorophenol-d4	7.767	132	49467	32.024	ng/ul	0.00
15) 4-Methylphenol-d8	8.941	113	53441	31.769	ng/ul	0.00
21) Nitrobenzene-d5	9.411	128	26643	32.485	ng/ul	0.00
24) 2-Nitrophenol-d4	10.140	143	28942	32.294	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.692	165	51730	29.608	ng/ul	0.00
31) 4-Chloroaniline-d4	11.203	131	72475	32.702	ng/ul	0.00
46) Dimethylphthalate-d6	14.270	166	192403	33.269	ng/ul	0.00
49) Acenaphthylene-d8	14.575	160	237862	32.997	ng/ul	0.00
54) 4-Nitrophenol-d4	15.068	143	24675m	28.171	ng/ul	0.00
60) Fluorene-d10	15.867	176	167267	33.013	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.979	200	26229	23.317	ng/ul	0.00
73) Anthracene-d10	17.724	188	297398	34.622	ng/ul	0.00
81) Pyrene-d10	20.003	212	387407	31.628	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.172	264	361349	34.653	ng/ul	0.00

Target Compounds Qvalue

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

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