

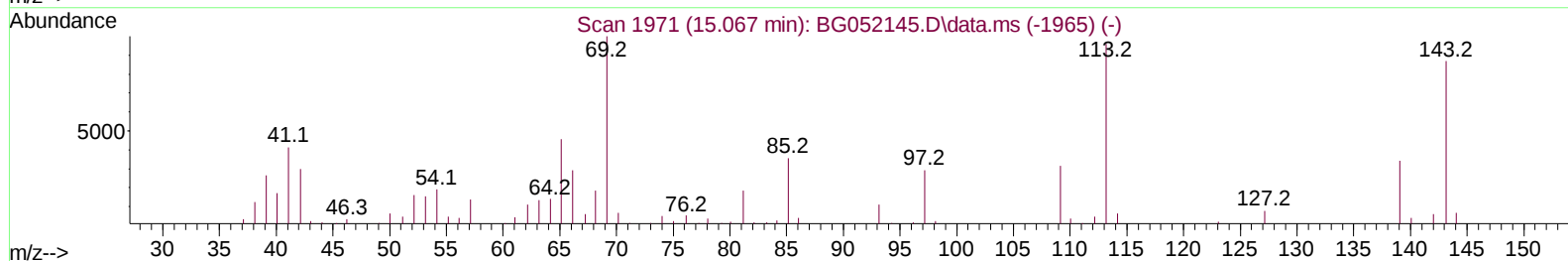
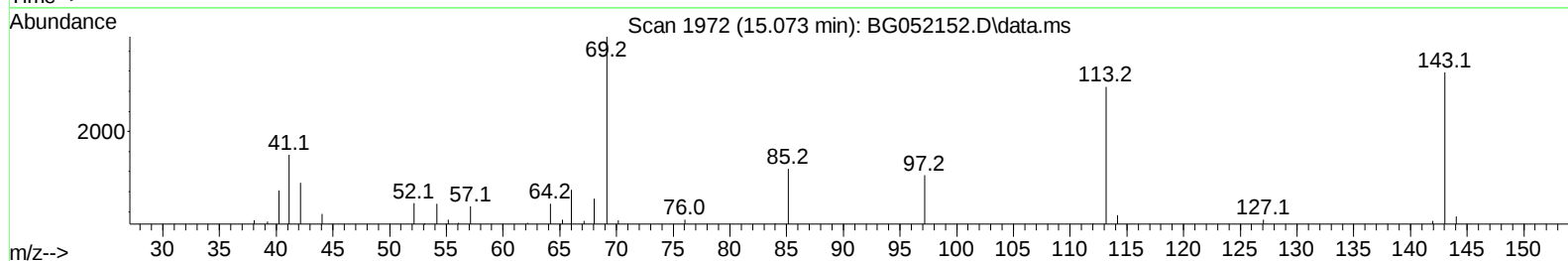
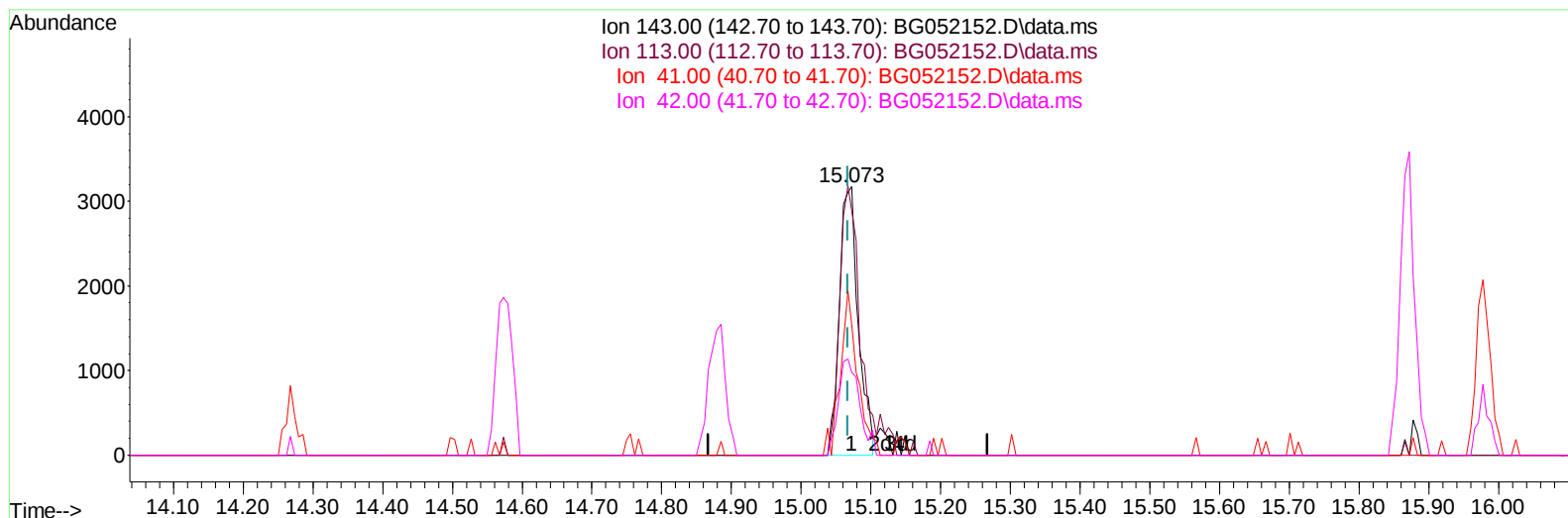
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052152.D
 Acq On : 26 Jan 2022 16:55
 Operator : CG/JU
 Sample : N1230-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0Q03

Manual IntegrationsAPPROVED

Quant Time: Jan 27 00:00:57 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 26 23:58:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/27/2022
 Supervised By :Yogesh Patel 01/28/2022



TIC: BG052152.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.073min (+ 0.005) 6.05 ng/u1

response 5920

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	91.15
41.00	44.40	48.50
42.00	29.70	30.93

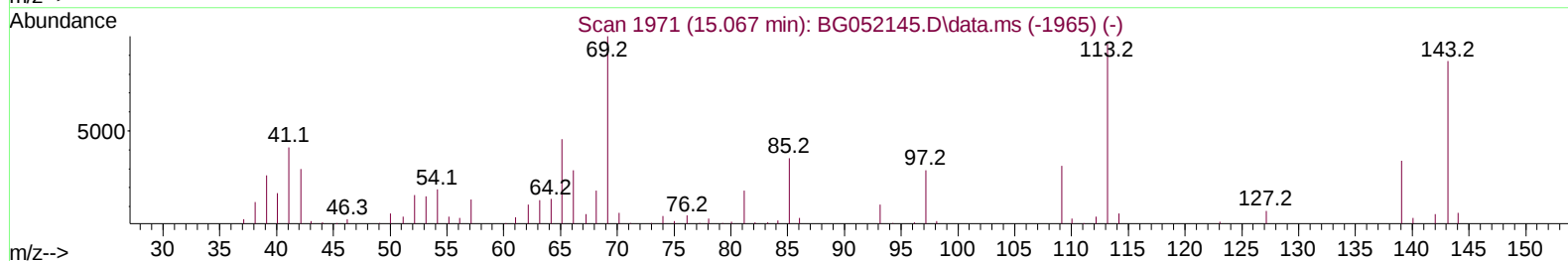
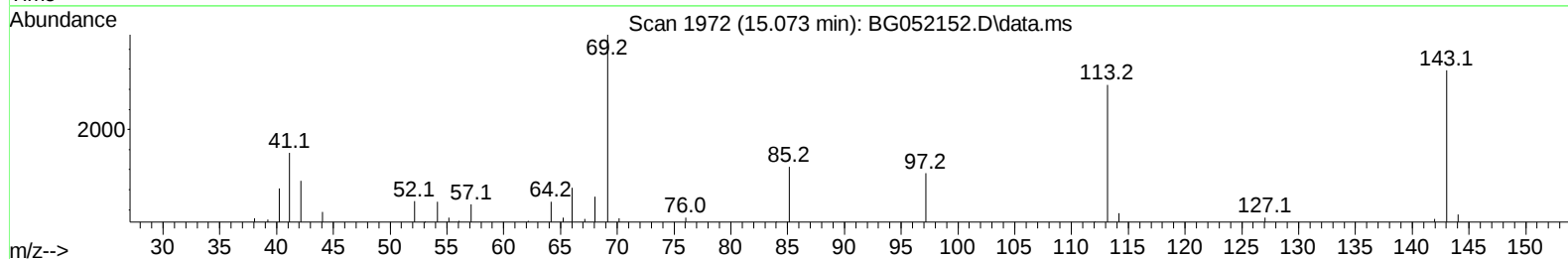
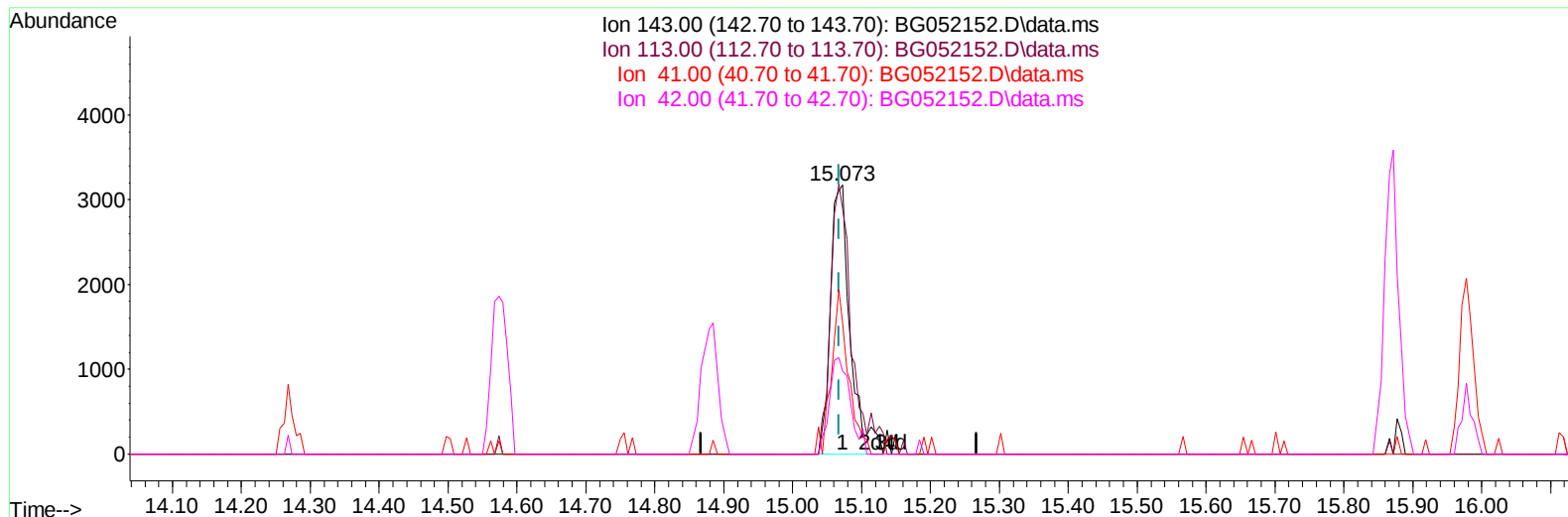
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TIC: BG052152.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.073min (+ 0.005) 6.41 ng/ul m

response 6278

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	91.15
41.00	44.40	48.50
42.00	29.70	30.93

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.241	152	24386	20.000	ng/ul	0.00
20) Naphthalene-d8	11.078	136	97742	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.885	164	70862	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.628	188	178055	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.946	240	194985	20.000	ng/ul	# 0.00
88) Perylene-d12	25.423	264	197435	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.565	96	2879	4.219	ng/uL	0.00
4) Pyridine-d5	3.994	84	19003	8.957	ng/ul	0.00
7) Phenol-d5	7.383	99	15870	6.649	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.548	67	44277	29.224	ng/ul	0.00
11) 2-Chlorophenol-d4	7.765	132	40375	25.550	ng/ul	0.00
15) 4-Methylphenol-d8	8.940	113	28996	15.628	ng/ul	0.00
21) Nitrobenzene-d5	9.410	128	27655	35.807	ng/ul	0.00
24) 2-Nitrophenol-d4	10.138	143	29040	33.998	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.690	165	50652	30.338	ng/ul	0.00
31) 4-Chloroaniline-d4	11.201	131	69361	29.831	ng/ul	0.00
46) Dimethylphthalate-d6	14.274	166	205345	35.058	ng/ul	0.00
49) Acenaphthylene-d8	14.579	160	245472	34.694	ng/ul	0.00
54) 4-Nitrophenol-d4	15.073	143	6278m	6.411	ng/ul	0.00
60) Fluorene-d10	15.872	176	183266	35.886	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.977	200	31790	28.302	ng/ul	0.00
73) Anthracene-d10	17.728	188	325557	38.543	ng/ul	0.00
81) Pyrene-d10	20.007	212	425776	37.942	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.188	264	402146	38.691	ng/ul	0.00

Target Compounds Qvalue

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

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