

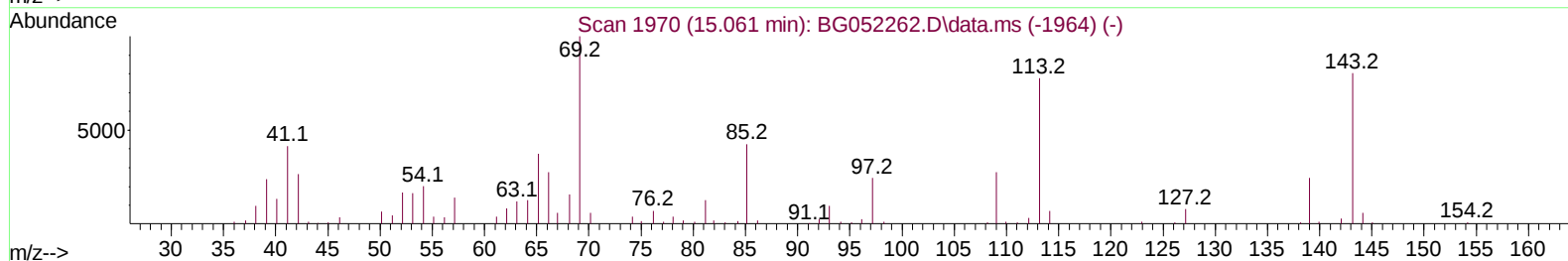
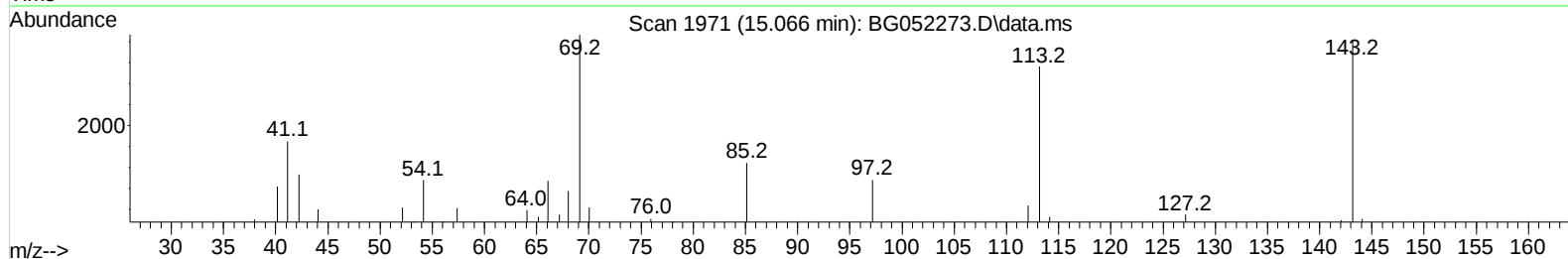
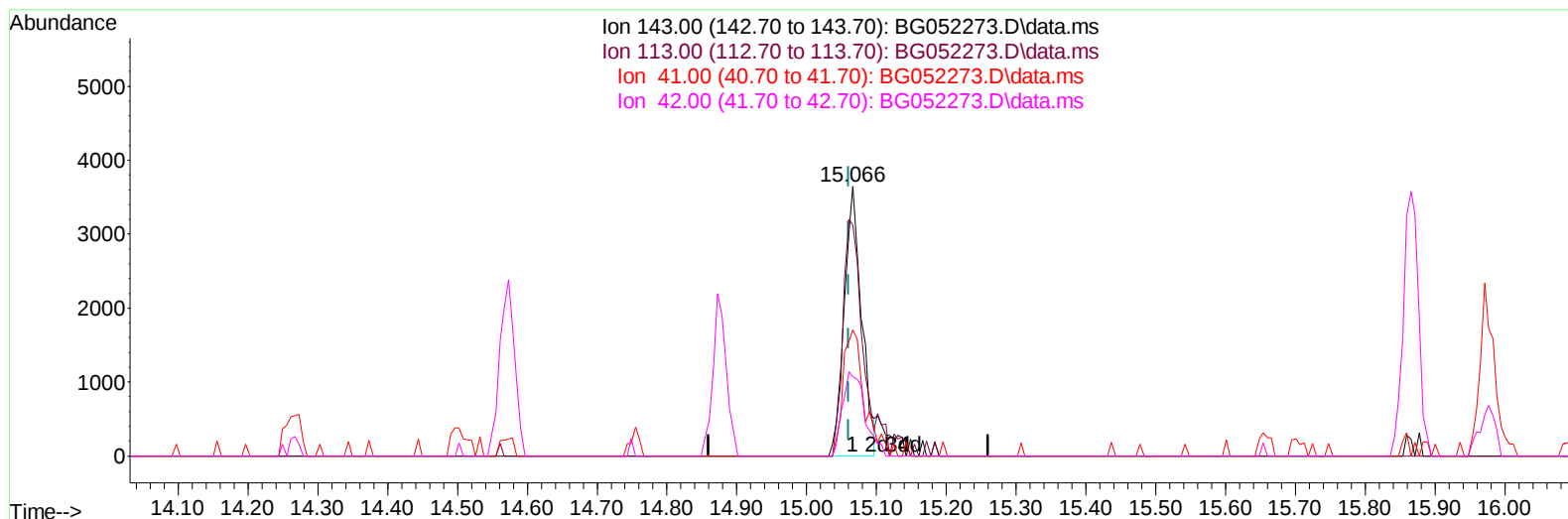
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020322\
Data File : BG052273.D
Acq On : 3 Feb 2022 17:48
Operator : CG/JU
Sample : N1349-12
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0HQ1

Manual IntegrationsAPPROVED

Quant Time: Feb 04 00:01:49 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 02/04/2022
Supervised By :Yogesh Patel 02/08/2022



TIC: BG052273.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.066min (+ 0.006) 6.41 ng/u1

response 6270

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	85.76
41.00	44.40	46.69
42.00	29.70	29.44

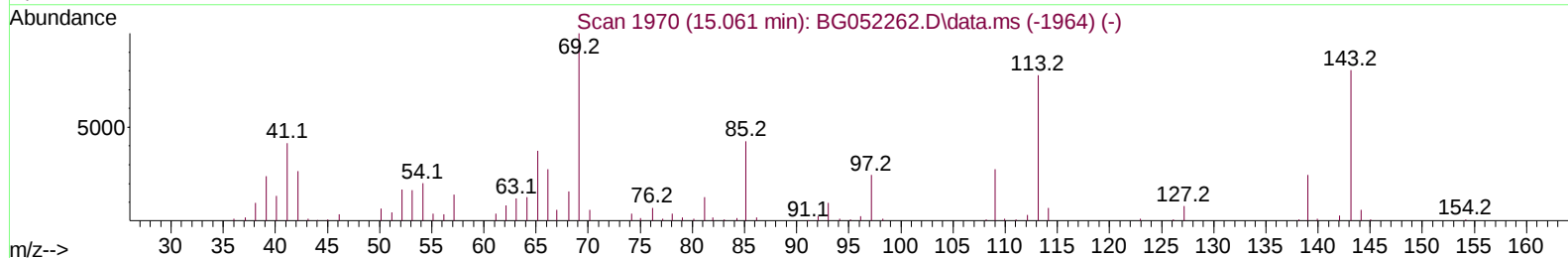
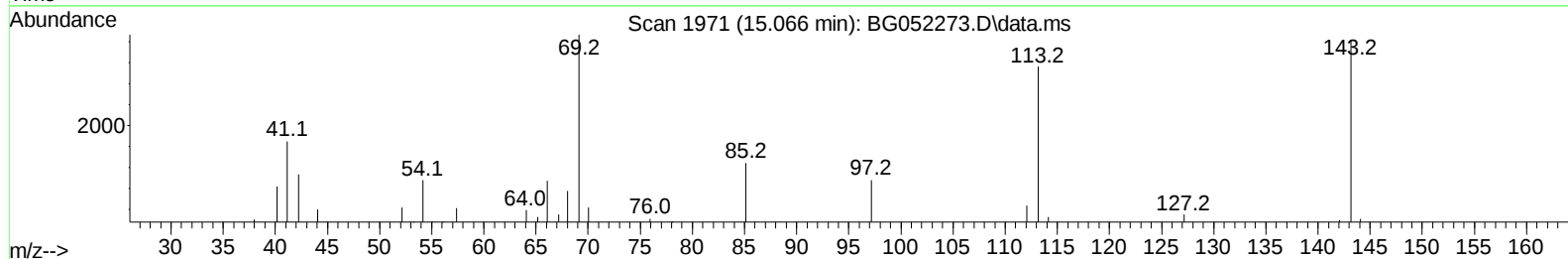
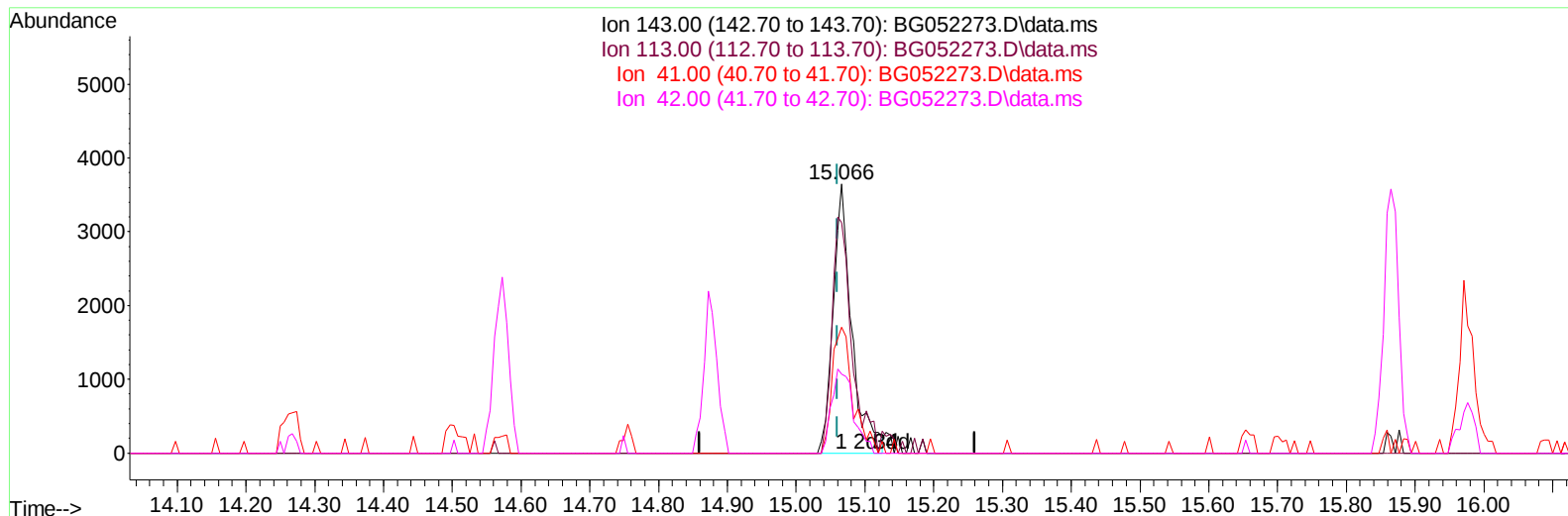
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TIC: BG052273.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.066min (+ 0.006) 7.03 ng/ul m

response 6879

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	85.76
41.00	44.40	46.69
42.00	29.70	29.44

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.234	152	29627	20.000	ng/ul	0.00
20) Naphthalene-d8	11.072	136	119885	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.878	164	82153	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.622	188	203020	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.939	240	218195	20.000	ng/ul	# 0.00
88) Perylene-d12	25.411	264	218612	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.564	96	3145	4.283	ng/uL	0.00
4) Pyridine-d5	4.005	84	6868	3.007	ng/ul	0.02
7) Phenol-d5	7.377	99	19141	7.103	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	56807	34.698	ng/ul	0.00
11) 2-Chlorophenol-d4	7.764	132	48515	25.835	ng/ul	0.00
15) 4-Methylphenol-d8	8.939	113	32394	15.840	ng/ul	0.00
21) Nitrobenzene-d5	9.403	128	33355	34.170	ng/ul	0.00
24) 2-Nitrophenol-d4	10.138	143	35058	32.867	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.684	165	60701	29.191	ng/ul	0.00
31) 4-Chloroaniline-d4	11.195	131	49833	18.893	ng/ul	0.00
46) Dimethylphthalate-d6	14.267	166	239847	37.142	ng/ul	0.00
49) Acenaphthylene-d8	14.573	160	290572	36.101	ng/ul	0.00
54) 4-Nitrophenol-d4	15.066	143	6879m	7.034	ng/ul	0.00
60) Fluorene-d10	15.865	176	214160	37.855	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.977	200	39771	31.731	ng/ul	0.00
73) Anthracene-d10	17.727	188	382126	39.925	ng/ul	0.00
81) Pyrene-d10	20.001	212	479294	36.652	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.176	264	461527	39.739	ng/ul	0.00

Target Compounds Qvalue

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

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