

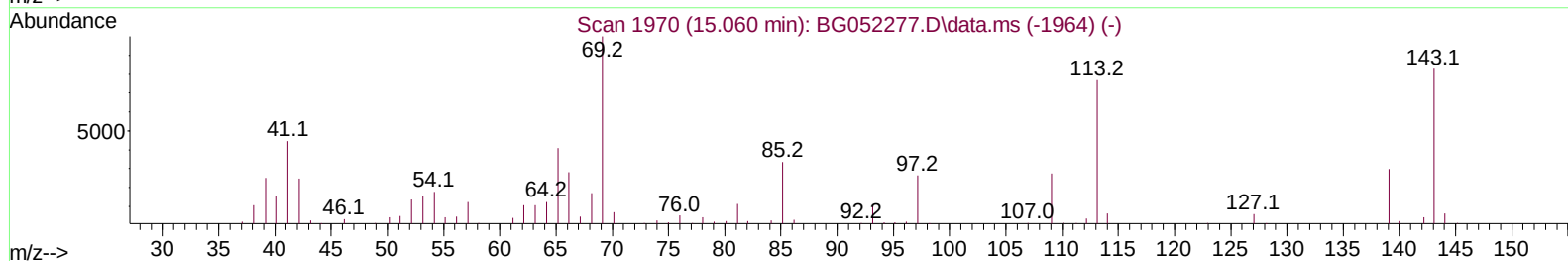
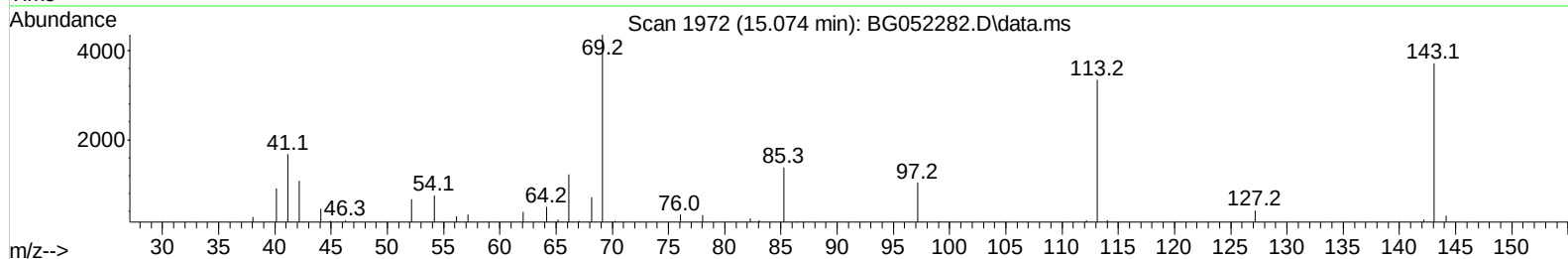
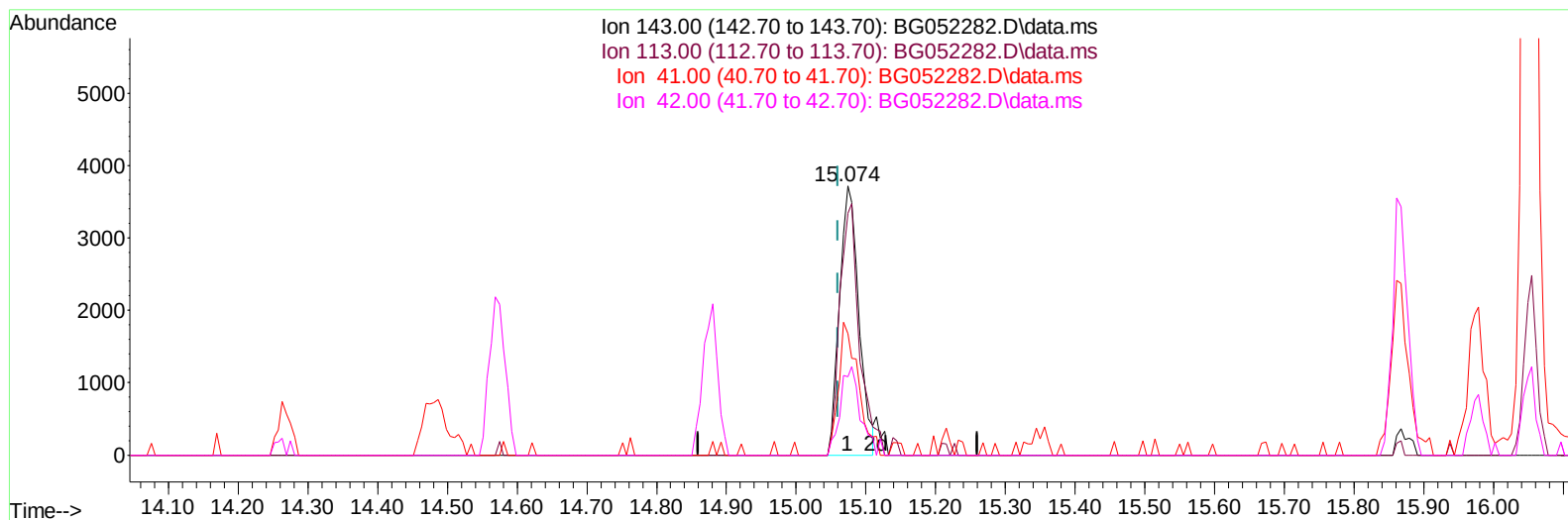
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020322\
Data File : BG052282.D
Acq On : 4 Feb 2022 00:06
Operator : CG/JU
Sample : N1328-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COHN2

Manual IntegrationsAPPROVED

Quant Time: Feb 04 01:38:08 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: BG052282.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.074min (+ 0.014) 6.72 ng/ul

response 7102

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	90.04
41.00	44.40	45.06
42.00	29.70	29.02

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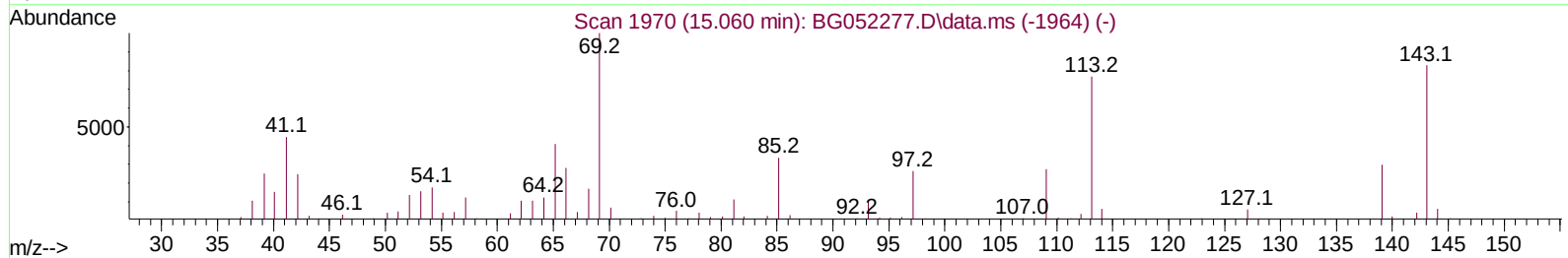
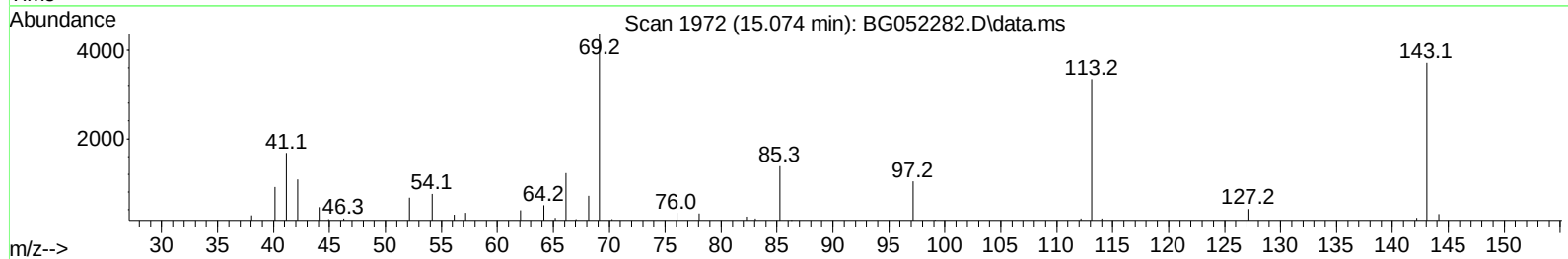
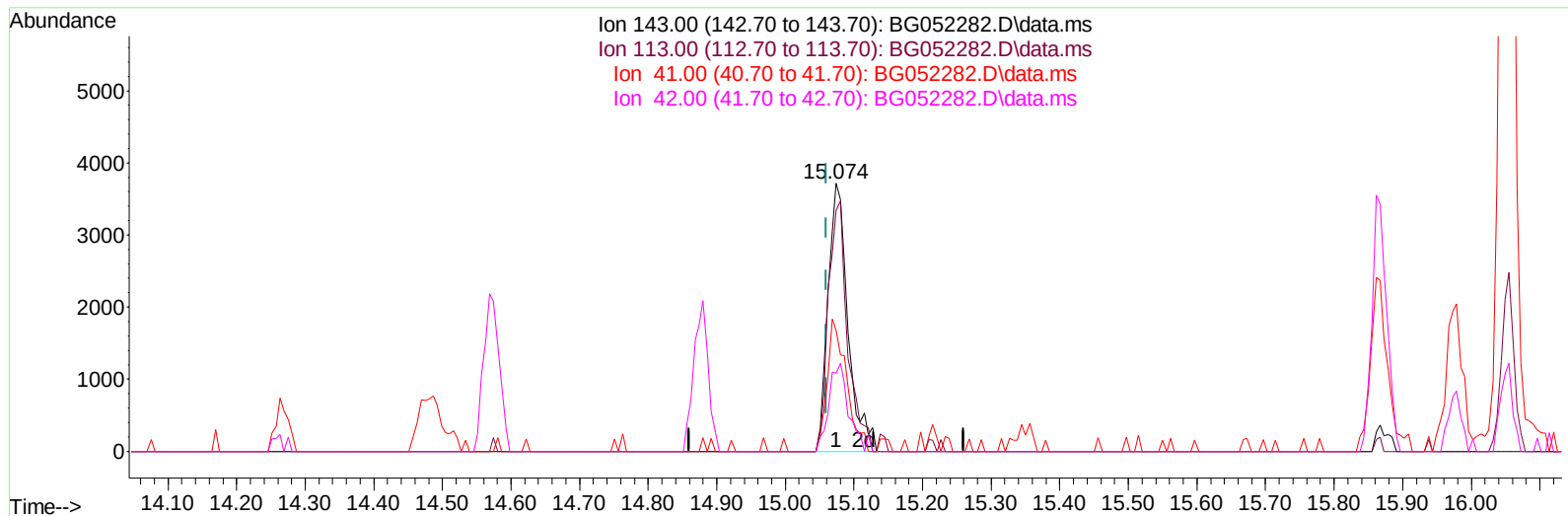
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TIC: BG052282.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.074min (+ 0.014) 7.08 ng/ul m

response 7485

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	90.04
41.00	44.40	45.06
42.00	29.70	29.02

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.236	152	31427	20.000	ng/ul	0.00
20) Naphthalene-d8	11.074	136	127262	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.880	164	88762	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.629	188	218560	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.941	240	227696	20.000	ng/ul	# 0.00
88) Perylene-d12	25.419	264	229099	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.566	96	3140	4.031	ng/uL	0.00
4) Pyridine-d5	3.995	84	10547	4.353	ng/ul	0.00
7) Phenol-d5	7.379	99	17620	6.164	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.537	67	53270	30.674	ng/ul	0.00
11) 2-Chlorophenol-d4	7.760	132	46461	23.324	ng/ul	0.00
15) 4-Methylphenol-d8	8.935	113	31939	14.723	ng/ul	0.00
21) Nitrobenzene-d5	9.405	128	32089	30.968	ng/ul	0.00
24) 2-Nitrophenol-d4	10.140	143	32567	28.762	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.686	165	59457	26.935	ng/ul	0.00
31) 4-Chloroaniline-d4	11.197	131	65299	23.321	ng/ul	0.00
46) Dimethylphthalate-d6	14.269	166	227194	32.563	ng/ul	0.00
49) Acenaphthylene-d8	14.575	160	271903	31.266	ng/ul	0.00
54) 4-Nitrophenol-d4	15.074	143	7485m	7.083	ng/ul	0.01
60) Fluorene-d10	15.867	176	197529	32.316	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.979	200	37805	28.018	ng/ul	0.00
73) Anthracene-d10	17.729	188	354466	34.401	ng/ul	0.00
81) Pyrene-d10	20.003	212	459288	33.656	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.178	264	429444	35.284	ng/ul	0.00

Target Compounds Qvalue

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

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