

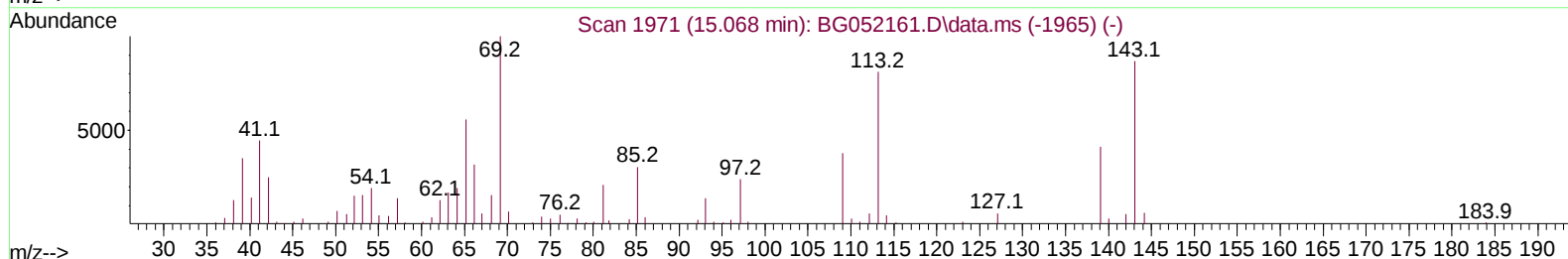
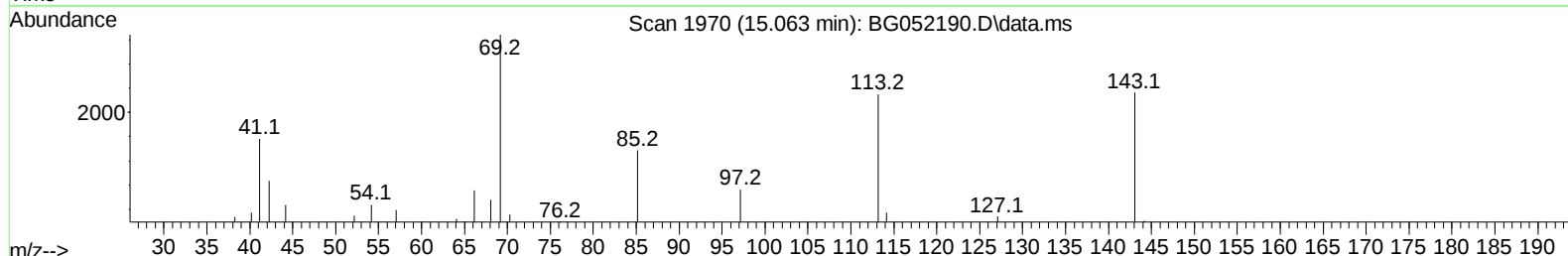
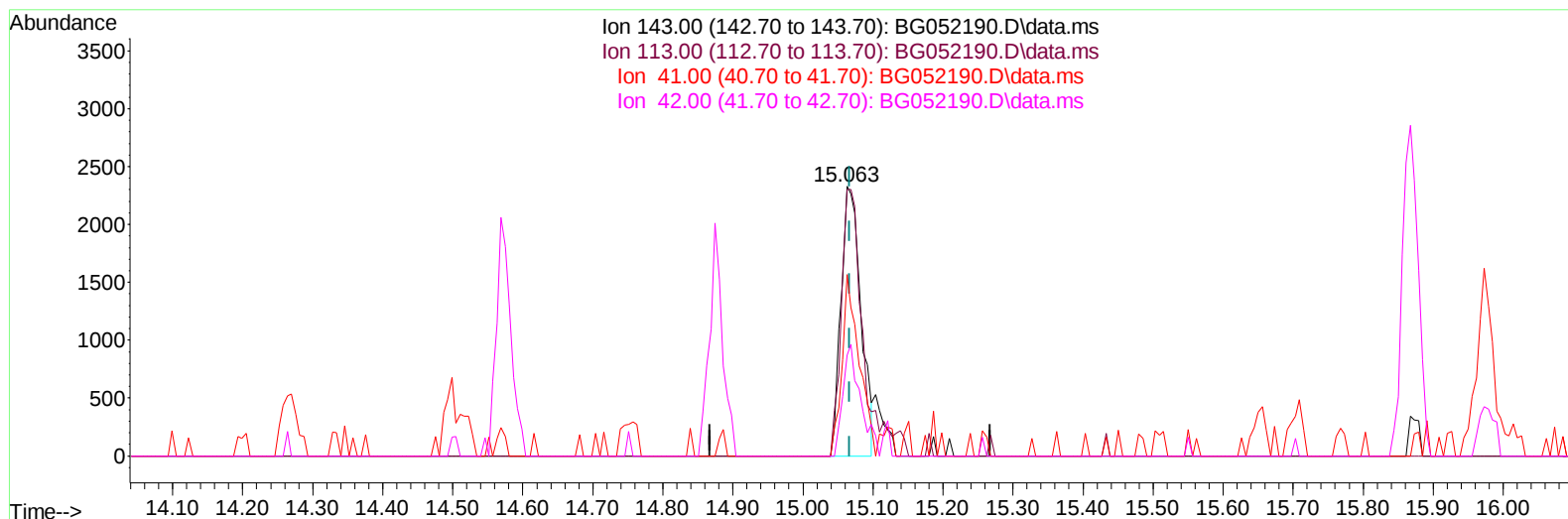
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
Data File : BG052190.D
Acq On : 27 Jan 2022 17:58
Operator : CG/JU
Sample : N1273-02
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0HK1

Manual IntegrationsAPPROVED

Quant Time: Jan 28 00:00:27 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 01/28/2022
Supervised By :Yogesh Patel 01/28/2022



TIC: BG052190.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.063min (-0.005) 4.70 ng/u1

response 4646

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	98.84#
41.00	44.40	67.23#
42.00	29.70	37.33#

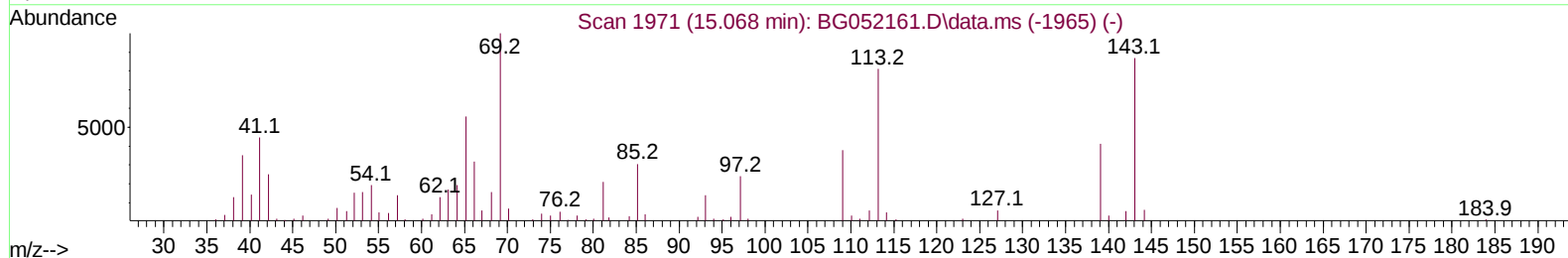
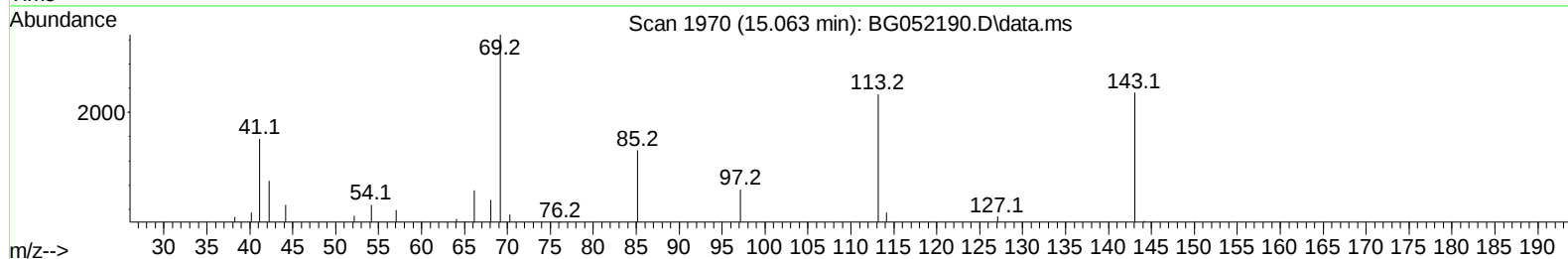
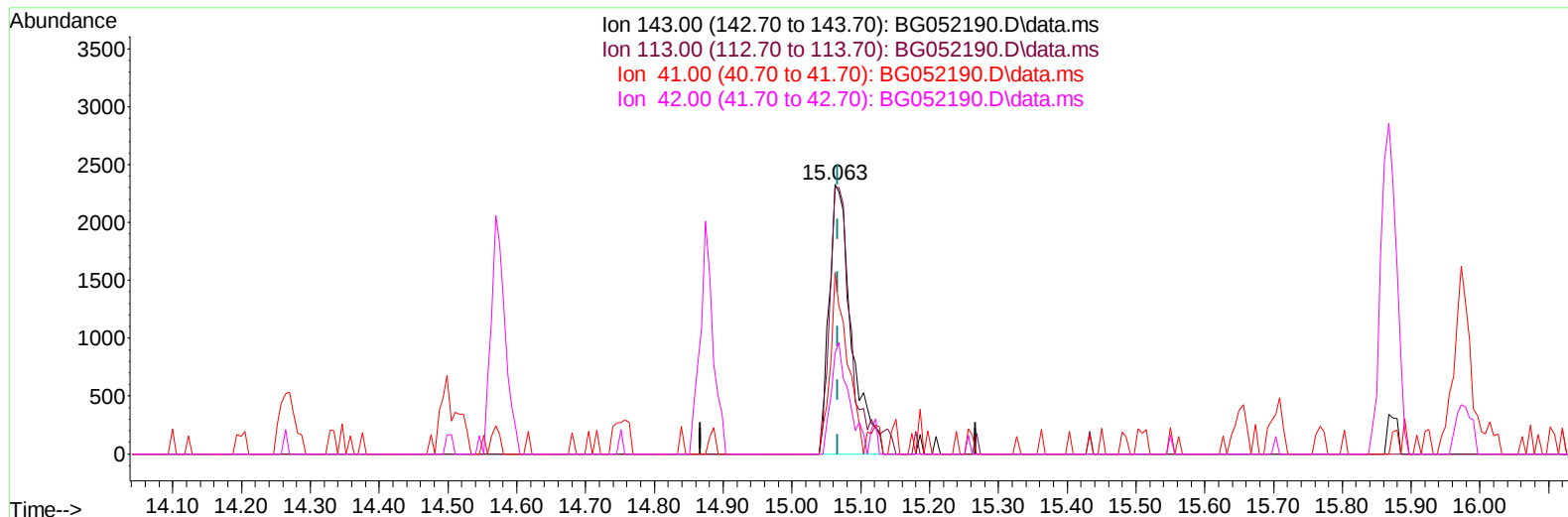
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(54) 4-Nitrophenol-d4 (S)

15.063min (-0.005) 5.27 ng/ul m

response 5213

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	98.84#
41.00	44.40	67.23#
42.00	29.70	37.33#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.237	152	22834	20.000	ng/ul	0.00
20) Naphthalene-d8	11.074	136	95564	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.880	164	71564	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.630	188	172829	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.942	240	176927	20.000	ng/ul	# 0.00
88) Perylene-d12	25.425	264	177821	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.566	96	2913	4.559	ng/uL	0.00
4) Pyridine-d5	3.995	84	13064	6.577	ng/ul	0.00
7) Phenol-d5	7.379	99	13573	6.073	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.543	67	39557	27.883	ng/ul	0.00
11) 2-Chlorophenol-d4	7.767	132	35595	24.056	ng/ul	0.00
15) 4-Methylphenol-d8	8.941	113	25200	14.505	ng/ul	0.00
21) Nitrobenzene-d5	9.406	128	25359	33.583	ng/ul	0.00
24) 2-Nitrophenol-d4	10.140	143	27316	32.708	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.686	165	46579	28.534	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.197	131	64005	28.155	ng/ul	-0.01
46) Dimethylphthalate-d6	14.270	166	199797	33.776	ng/ul	0.00
49) Acenaphthylene-d8	14.575	160	237386	33.222	ng/ul	0.00
54) 4-Nitrophenol-d4	15.063	143	5213m	5.271	ng/ul	0.00
60) Fluorene-d10	15.867	176	177435	34.404	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.979	200	25886	23.742	ng/ul	0.00
73) Anthracene-d10	17.730	188	310607	37.885	ng/ul	0.00
81) Pyrene-d10	20.009	212	366621	36.005	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.184	264	344572	36.808	ng/ul	0.00

Target Compounds Qvalue

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

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