

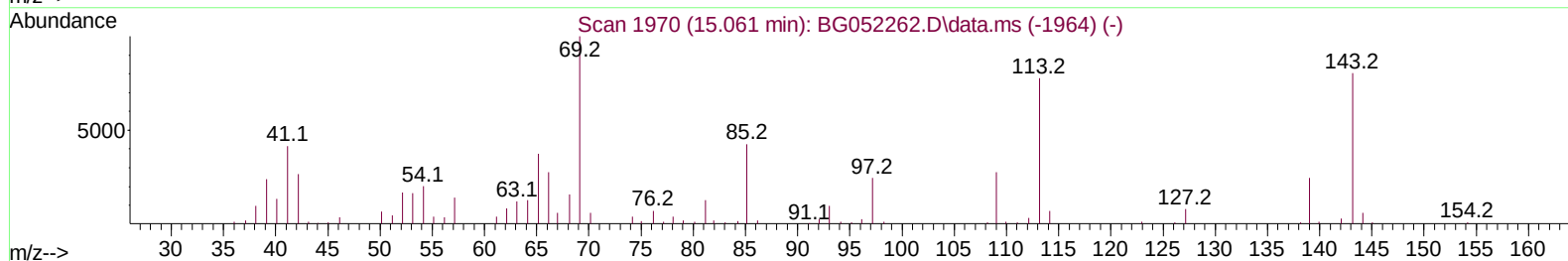
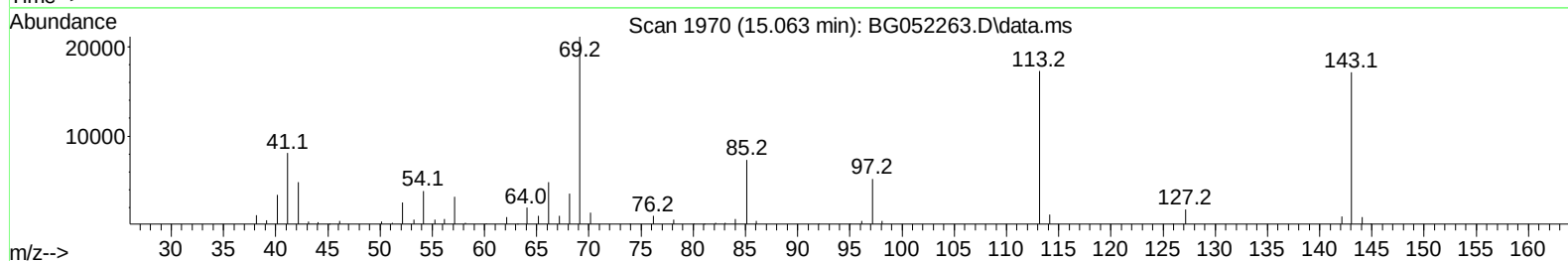
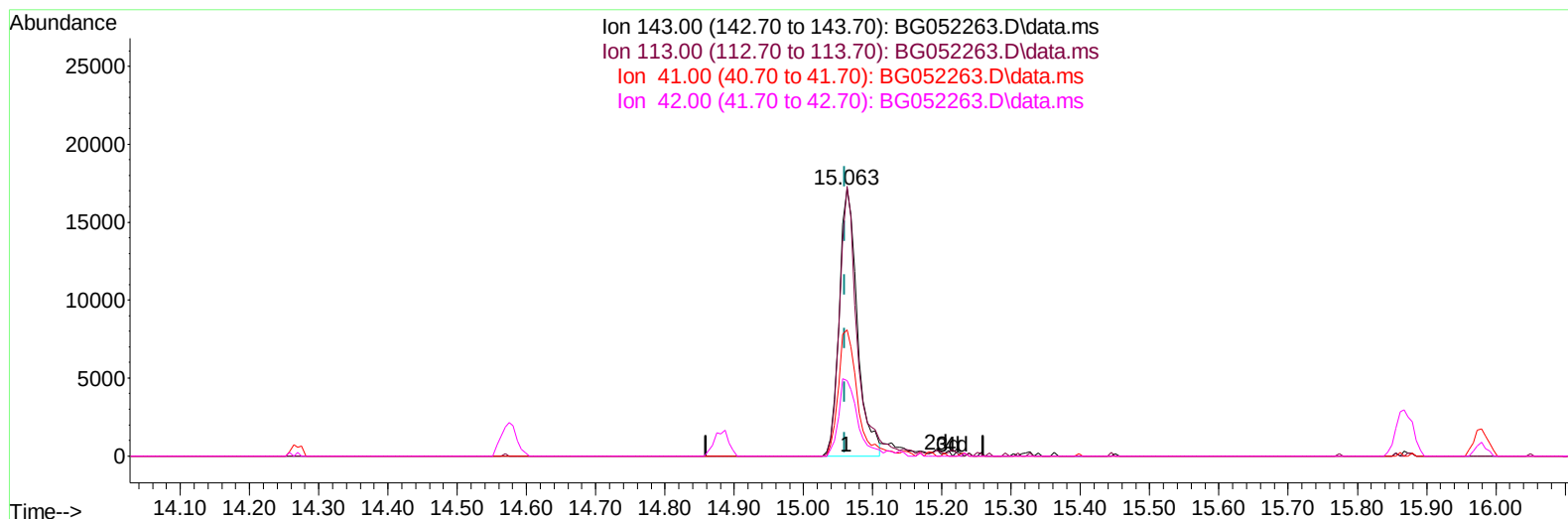
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020322\
Data File : BG052263.D
Acq On : 3 Feb 2022 11:15
Operator : CG/JU
Sample : PB142430BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK430

Manual IntegrationsAPPROVED

Quant Time: Feb 03 23:59:11 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: BG052263.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.063min (+ 0.003) 33.09 ng/ul

response 31153

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	100.98#
41.00	44.40	47.31
42.00	29.70	28.27

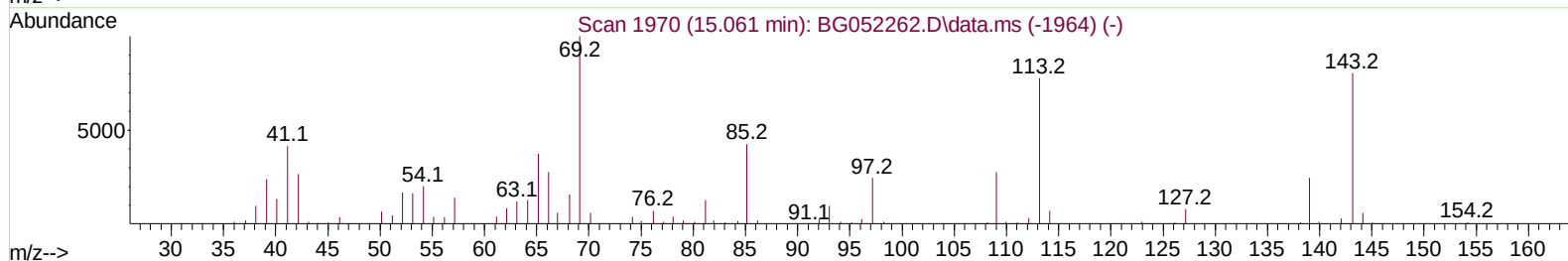
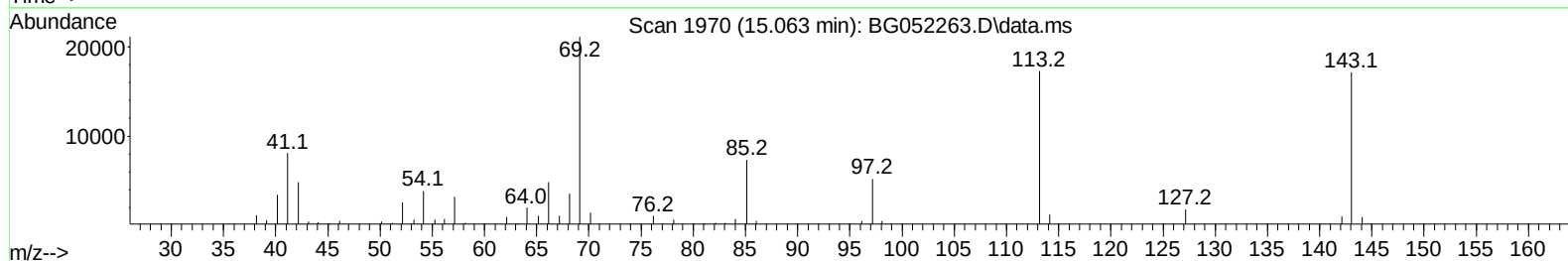
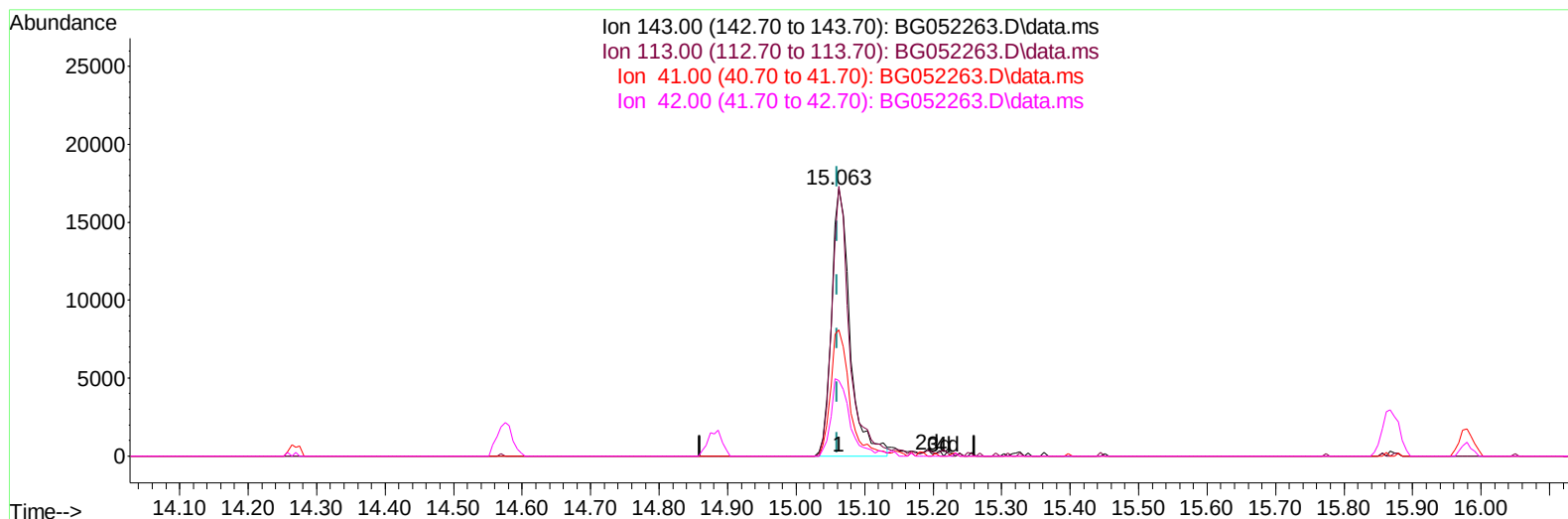
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TIC: BG052263.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.063min (+ 0.003) 34.20 ng/ul m

response 32200

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	100.98#
41.00	44.40	47.31
42.00	29.70	28.27

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.242	152	28270	20.000	ng/ul	0.00
20) Naphthalene-d8	11.080	136	112699	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.881	164	79085	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.630	188	192334	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.942	240	203053	20.000	ng/ul	# 0.00
88) Perylene-d12	25.413	264	200259	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.566	96	5073	7.239	ng/uL	0.00
4) Pyridine-d5	3.995	84	70153	32.187	ng/ul	0.00
7) Phenol-d5	7.385	99	82884	32.233	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.543	67	51401	32.903	ng/ul	0.00
11) 2-Chlorophenol-d4	7.767	132	59946	33.454	ng/ul	0.00
15) 4-Methylphenol-d8	8.947	113	63551	32.567	ng/ul	0.00
21) Nitrobenzene-d5	9.411	128	31038	33.824	ng/ul	0.00
24) 2-Nitrophenol-d4	10.146	143	34668	34.574	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.692	165	65646	33.582	ng/ul	0.00
31) 4-Chloroaniline-d4	11.203	131	85358	34.424	ng/ul	0.00
46) Dimethylphthalate-d6	14.270	166	225628	36.296	ng/ul	0.00
49) Acenaphthylene-d8	14.575	160	279563	36.081	ng/ul	0.00
54) 4-Nitrophenol-d4	15.063	143	32200m	34.201	ng/ul	0.00
60) Fluorene-d10	15.867	176	200339	36.786	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.979	200	34819	29.323	ng/ul	0.00
73) Anthracene-d10	17.730	188	343670	37.902	ng/ul	0.00
81) Pyrene-d10	20.003	212	432978	35.579	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.178	264	400287	37.625	ng/ul	0.00

Target Compounds Qvalue

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

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