

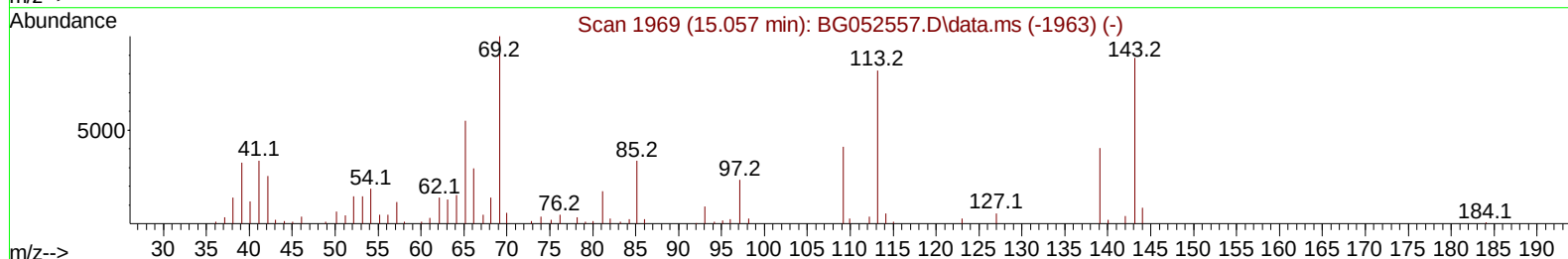
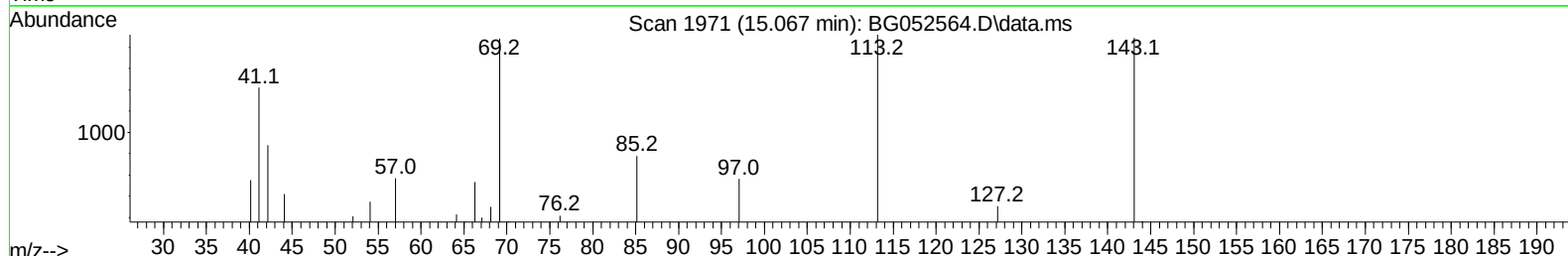
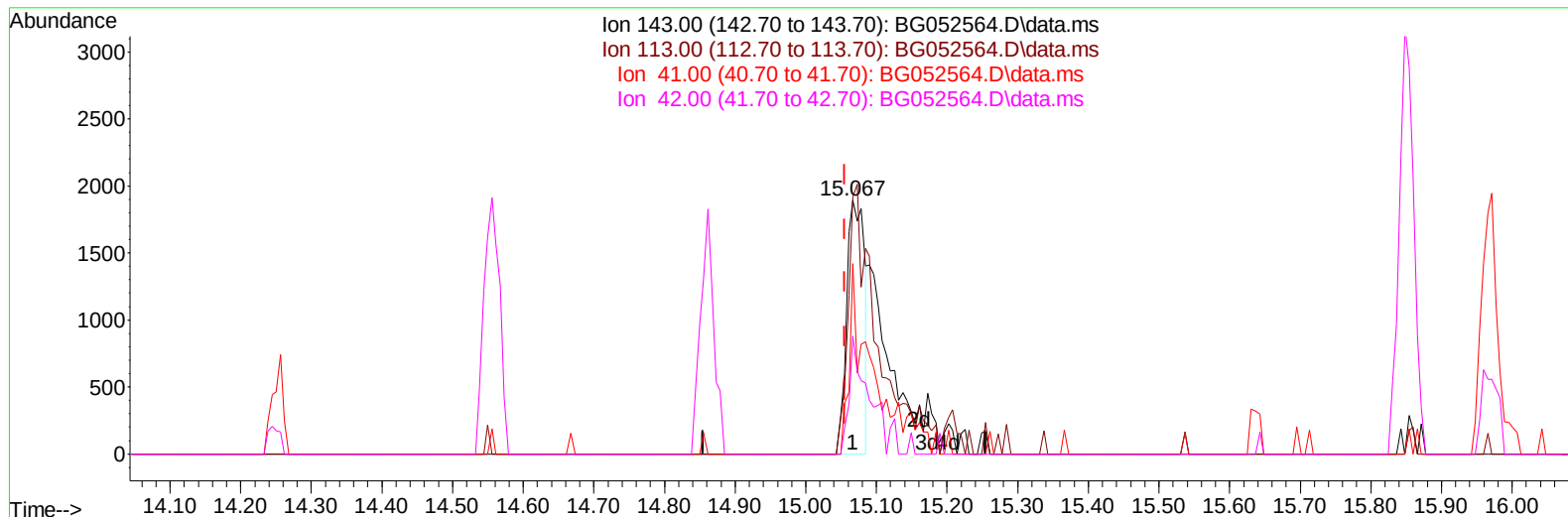
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030322\
Data File : BG052564.D
Acq On : 3 Mar 2022 21:14
Operator : CG/JU
Sample : N1745-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
H0AB1

Manual IntegrationsAPPROVED

Quant Time: Mar 04 00:03:35 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022222.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 22 17:32:12 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/04/2022
Supervised By :Yogesh Patel 03/07/2022



TIC: BG052564.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.067min (+ 0.012) 3.24 ng/ul

response 3315

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	101.43#
41.00	44.40	75.11#
42.00	29.70	46.41#

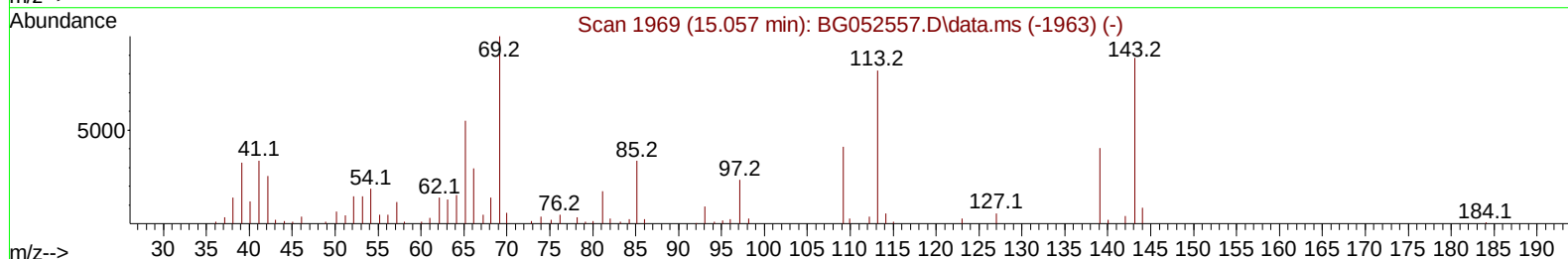
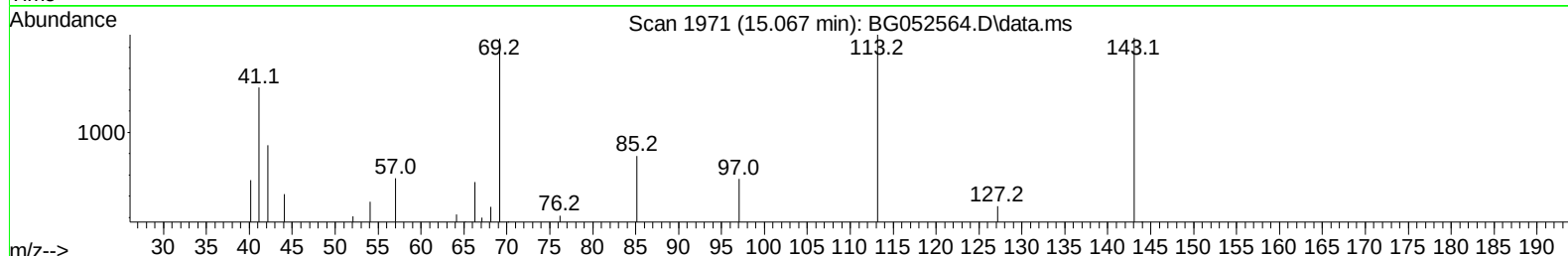
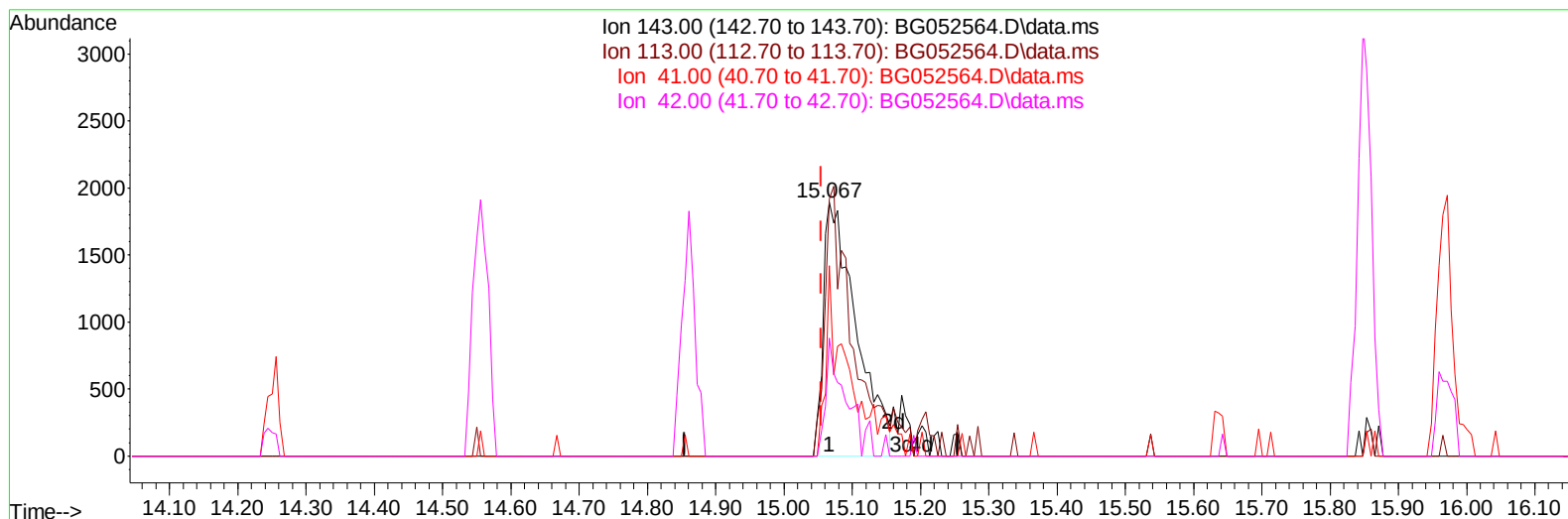
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TIC: BG052564.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.067min (+ 0.012) 6.16 ng/ul m

response 6311

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	101.43#
41.00	44.40	75.11#
42.00	29.70	46.41#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.217	152	24927	20.000	ng/ul	0.00
20) Naphthalene-d8	11.055	136	104205	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.861	164	79278	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.610	188	205711	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.928	240	219318	20.000	ng/ul	# 0.00
88) Perylene-d12	25.388	264	216262	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.541	96	2698	4.246	ng/uL	0.00
4) Pyridine-d5	3.982	84	13360	6.969	ng/ul	0.00
7) Phenol-d5	7.365	99	17238	7.293	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.518	67	42847	31.018	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.747	132	39962	24.071	ng/ul	0.00
15) 4-Methylphenol-d8	8.922	113	29759	16.181	ng/ul	-0.01
21) Nitrobenzene-d5	9.386	128	28834	32.458	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.121	143	29936	30.765	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.673	165	51485	28.374	ng/ul	0.00
31) 4-Chloroaniline-d4	11.184	131	68101	27.082	ng/ul	0.00
46) Dimethylphthalate-d6	14.250	166	230800	35.812	ng/ul	0.00
49) Acenaphthylene-d8	14.556	160	273698	34.402	ng/ul	0.00
54) 4-Nitrophenol-d4	15.067	143	6311m	6.160	ng/ul	0.01
60) Fluorene-d10	15.854	176	203624	35.966	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.966	200	36126	27.697	ng/ul	0.00
73) Anthracene-d10	17.710	188	371828	37.415	ng/ul	0.00
81) Pyrene-d10	19.990	212	472477	36.816	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.147	264	453301	39.070	ng/ul	-0.01

Target Compounds Qvalue

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

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