

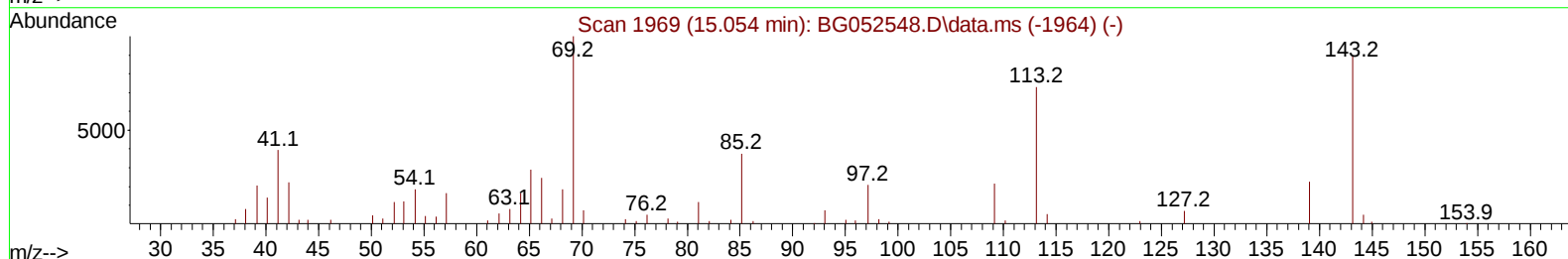
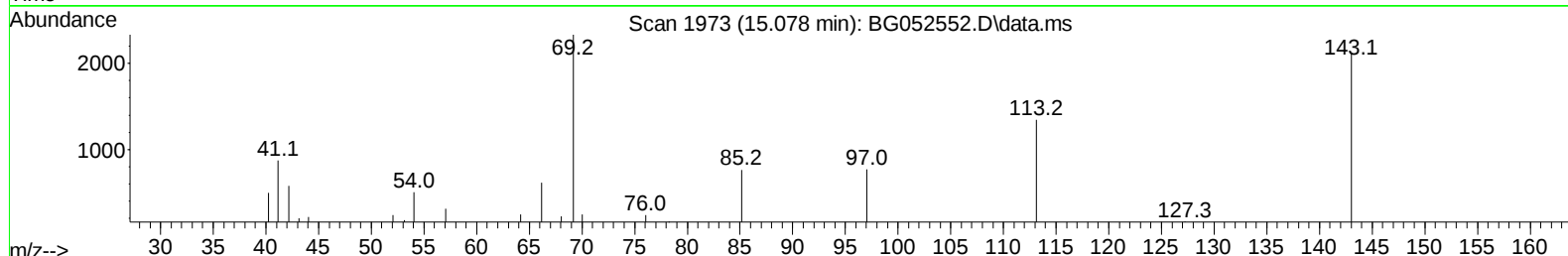
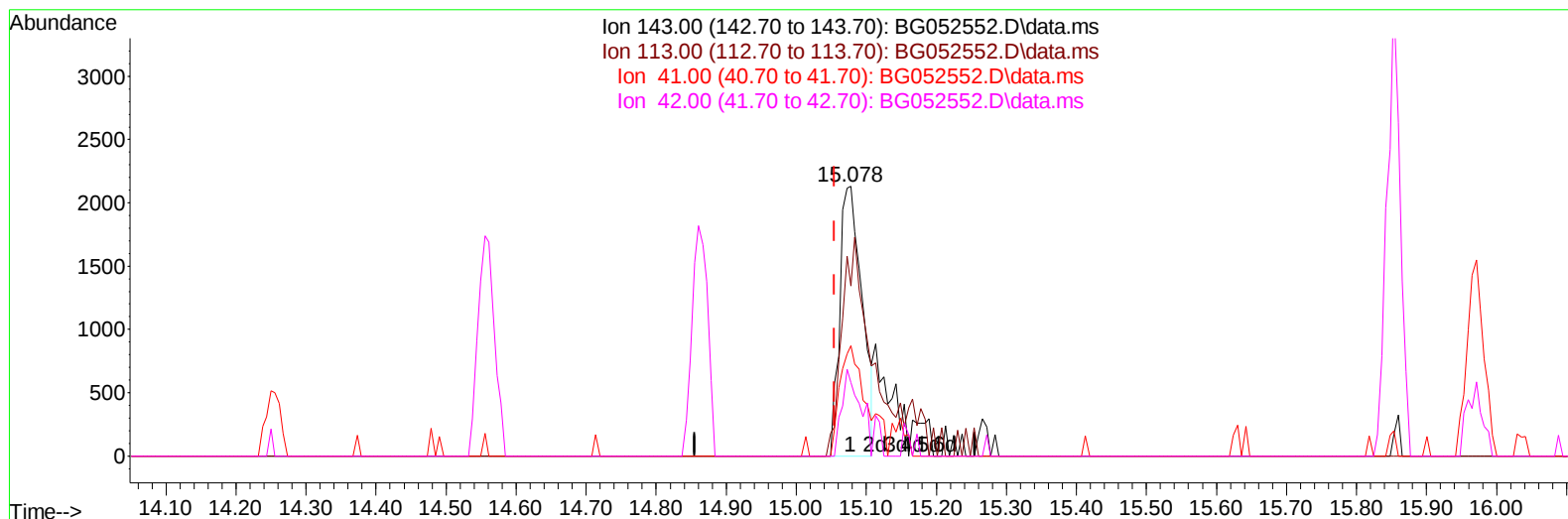
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030322\
Data File : BG052552.D
Acq On : 3 Mar 2022 13:31
Operator : CG/JU
Sample : N1745-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
H0AB0

Manual IntegrationsAPPROVED

Quant Time: Mar 04 00:00:57 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: BG052552.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.078min (+ 0.023) 4.00 ng/u1

response 4775

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	63.13#
41.00	44.40	40.96
42.00	29.70	27.20

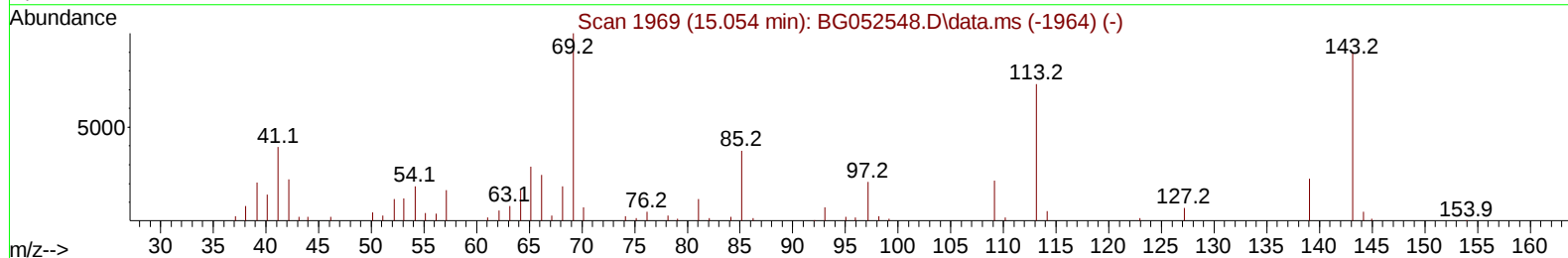
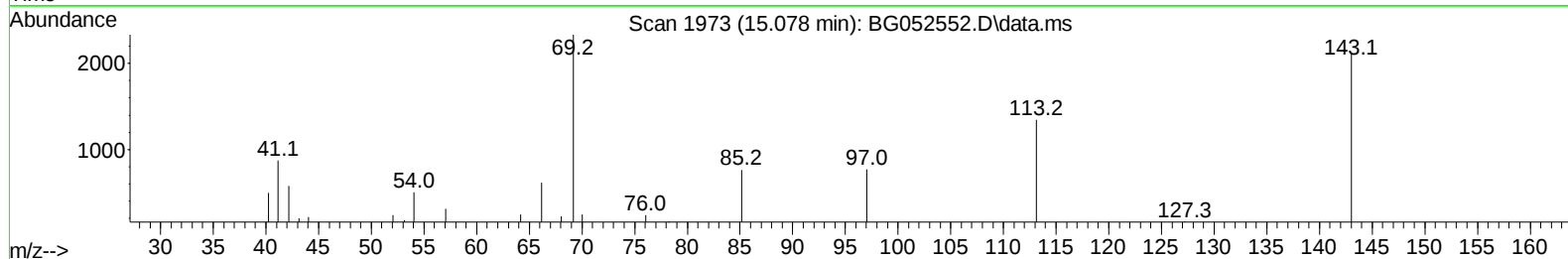
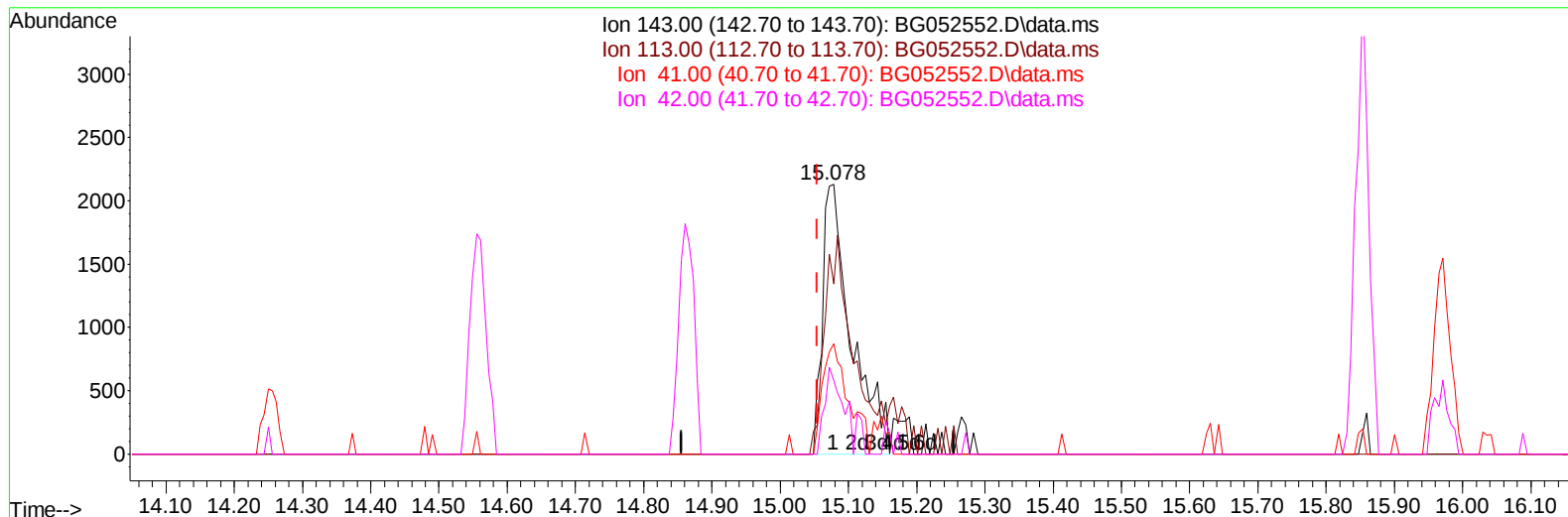
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TIC: BG052552.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.078min (+ 0.023) 5.23 ng/ul m

response 6234

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	63.13#
41.00	44.40	40.96
42.00	29.70	27.20

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.217	152	28499	20.000	ng/ul	0.00
20) Naphthalene-d8	11.054	136	121473	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.867	164	92312	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.616	188	230118	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.928	240	234434	20.000	ng/ul	# 0.00
88) Perylene-d12	25.393	264	237946	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.541	96	2733	3.762	ng/uL	0.00
4) Pyridine-d5	3.981	84	12677	5.784	ng/ul	0.00
7) Phenol-d5	7.371	99	17667	6.538	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.524	67	47023	29.775	ng/ul	0.00
11) 2-Chlorophenol-d4	7.747	132	43638	22.990	ng/ul	0.00
15) 4-Methylphenol-d8	8.928	113	31467	14.966	ng/ul	0.00
21) Nitrobenzene-d5	9.392	128	30263	29.224	ng/ul	0.00
24) 2-Nitrophenol-d4	10.126	143	32713	28.840	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.672	165	54884	25.947	ng/ul	0.00
31) 4-Chloroaniline-d4	11.183	131	68670	23.427	ng/ul	0.00
46) Dimethylphthalate-d6	14.256	166	241093	32.127	ng/ul	0.00
49) Acenaphthylene-d8	14.561	160	286428	30.919	ng/ul	0.00
54) 4-Nitrophenol-d4	15.078	143	6234m	5.226	ng/ul	0.02
60) Fluorene-d10	15.853	176	212903	32.295	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.971	200	37770	25.886	ng/ul	0.00
73) Anthracene-d10	17.716	188	387818	34.885	ng/ul	0.00
81) Pyrene-d10	19.995	212	477795	34.830	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.153	264	440150	34.479	ng/ul	0.00

Target Compounds Qvalue

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

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