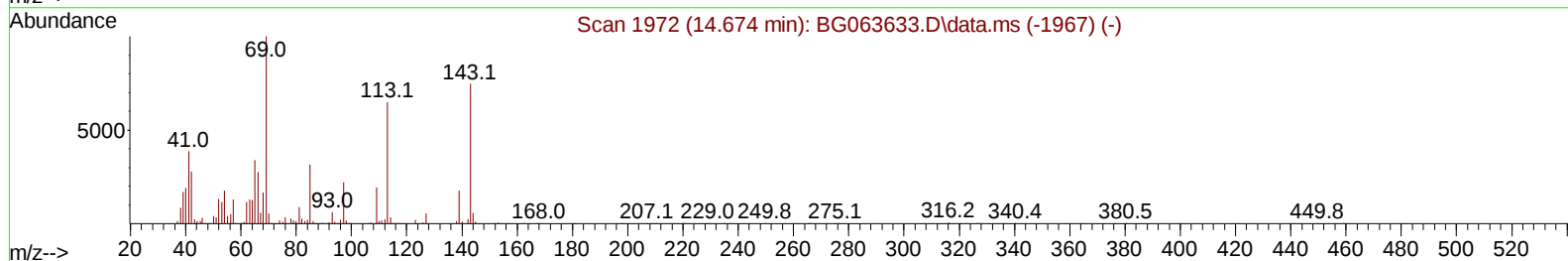
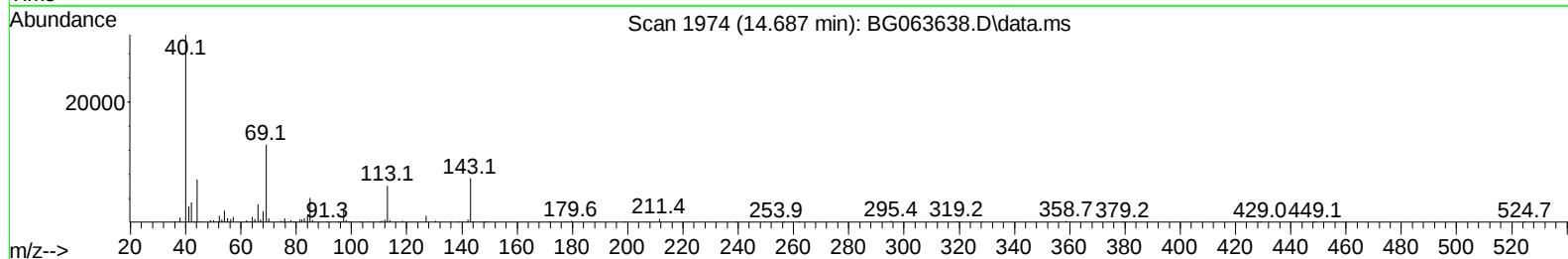
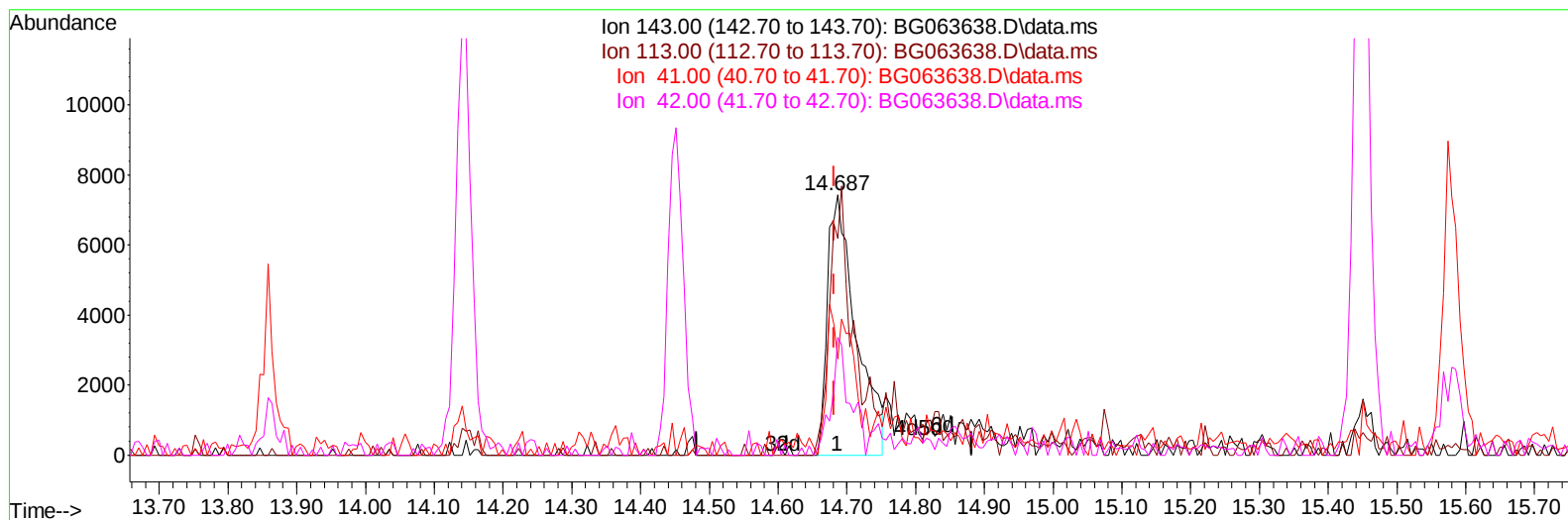


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
Data File : BG063638.D
Acq On : 10 Dec 2024 19:12
Operator : RC/JU
Sample : P5109-11
Misc :
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Dec 10 23:26:51 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 06 15:14:46 2024
Response via : Initial Calibration



TIC: BG063638.D\data.ms

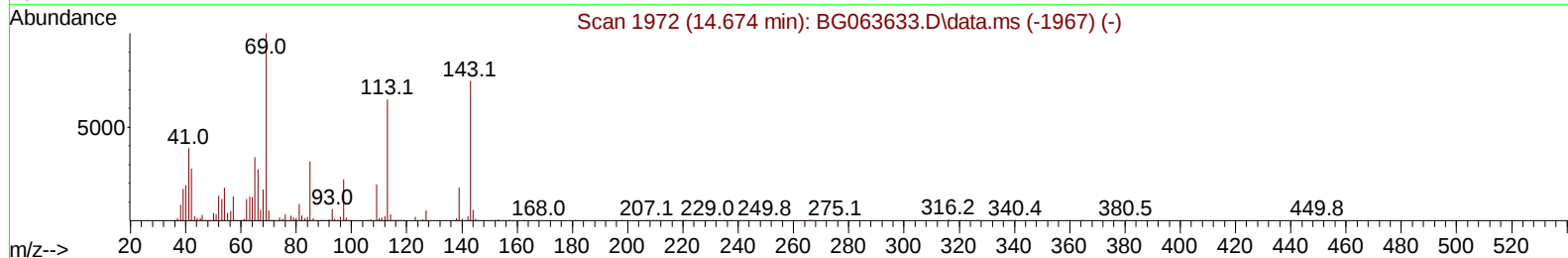
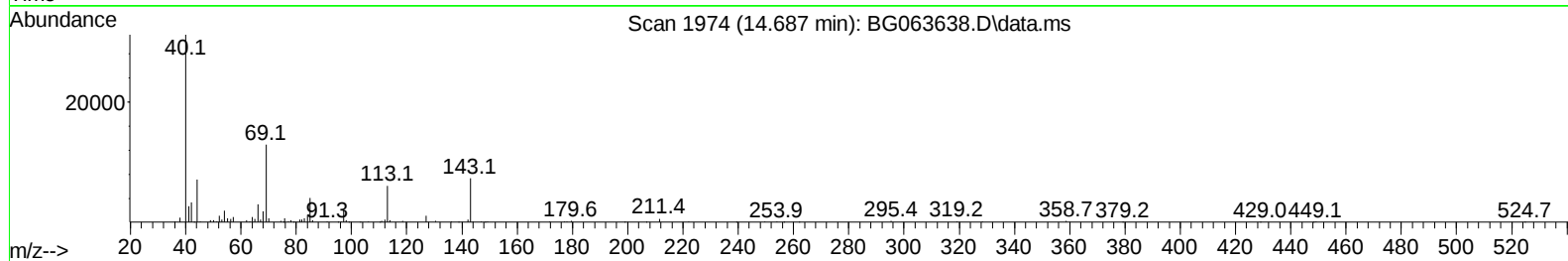
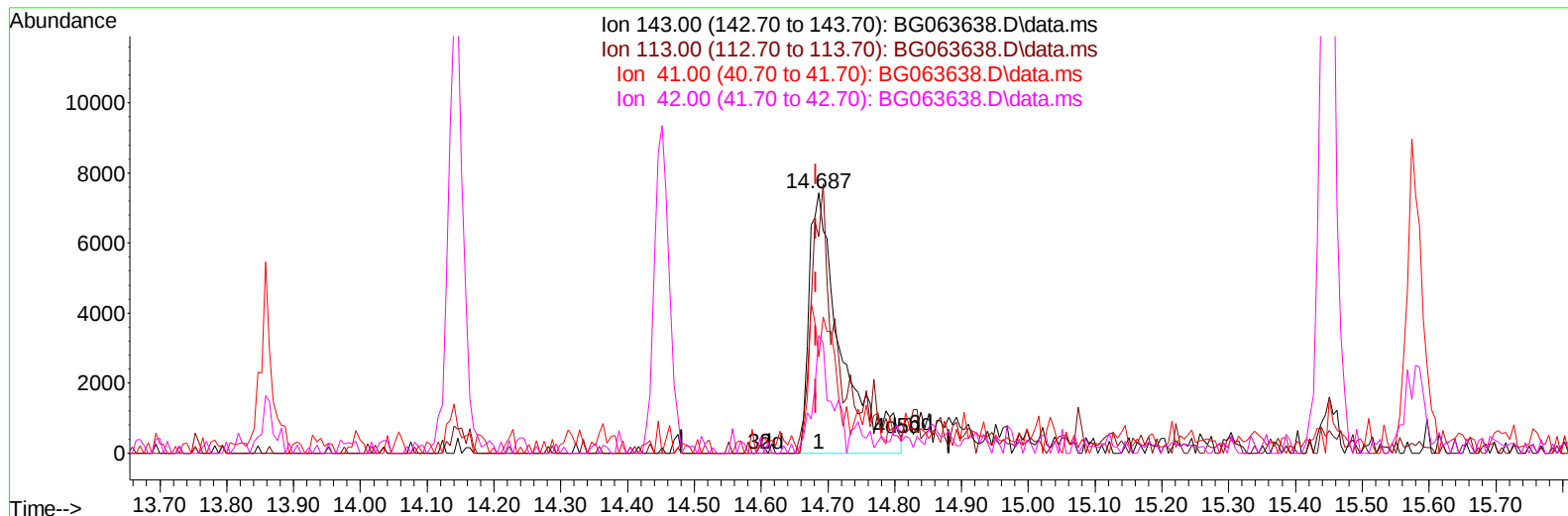
(54) 4-Nitrophenol-d4 (S)

14.687min (+ 0.005) 5.19 ng/ul

response 21358

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	110.90	83.31#
41.00	45.50	36.84
42.00	30.70	45.34#

Quant Time: Dec 10 23:26:51 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 06 15:14:46 2024
Response via : Initial Calibration



TIC: BG063638.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.687min (+ 0.005) 6.05 ng/u1 m

response 24884

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	110.90	83.31#
41.00	45.50	36.84
42.00	30.70	45.34#

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
 Data File : BG063638.D
 Acq On : 10 Dec 2024 19:12
 Operator : RC/JU
 Sample : P5109-11
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Dec 10 23:41:28 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 06 15:14:46 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.812	152	140570	20.000	ng/ul	0.00
20) Naphthalene-d8	10.603	136	584381	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.452	164	442215	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.201	188	1017637	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.455	240	958272	20.000	ng/ul	#-0.02
88) Perylene-d12	24.475	264	1044502	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.282	96	19239	6.042	ng/uL	0.00
4) Pyridine-d5	3.694	84	69808	5.926	ng/ul	0.00
7) Phenol-d5	6.996	99	124479	8.994	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.143	67	347460	37.130	ng/ul	0.00
11) 2-Chlorophenol-d4	7.348	132	287193	35.226	ng/ul	0.00
15) 4-Methylphenol-d8	8.529	113	232942	22.701	ng/ul	0.00
21) Nitrobenzene-d5	8.970	128	167635	40.868	ng/ul	0.00
24) 2-Nitrophenol-d4	9.687	143	187563	39.958	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.233	165	362790	35.896	ng/ul	0.00
31) 4-Chloroaniline-d4	10.738	131	334989	27.427	ng/ul	0.00
46) Dimethylphthalate-d6	13.858	166	1308448	40.712	ng/ul	0.00
49) Acenaphthylene-d8	14.146	160	1410549	40.152	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	24884m	6.047	ng/ul	0.00
60) Fluorene-d10	15.450	176	1152884	36.391	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.580	200	152016	26.443	ng/ul	0.00
73) Anthracene-d10	17.307	188	1985120	42.721	ng/ul	0.00
81) Pyrene-d10	19.604	212	2289834	43.110	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.275	264	2256544	43.014	ng/ul	-0.02

Target Compounds	Qvalue

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

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