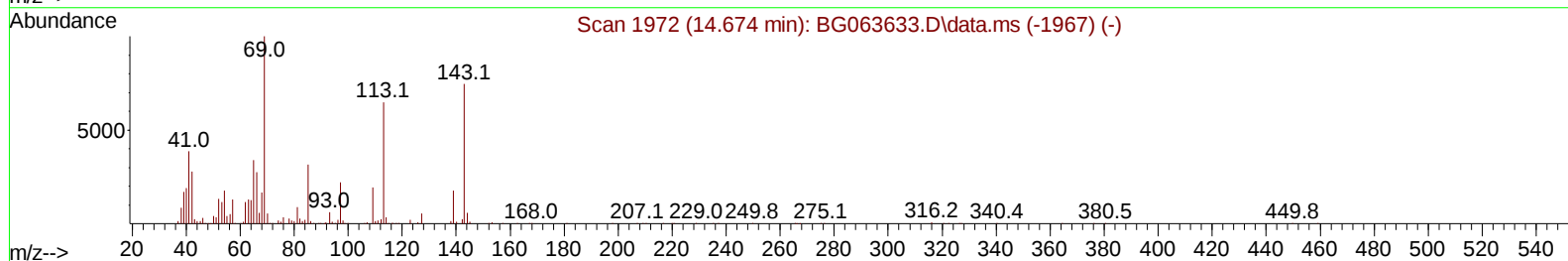
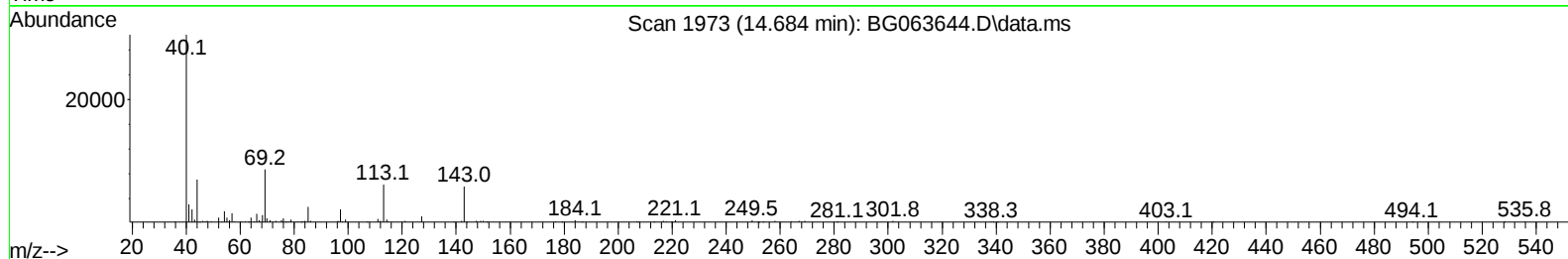
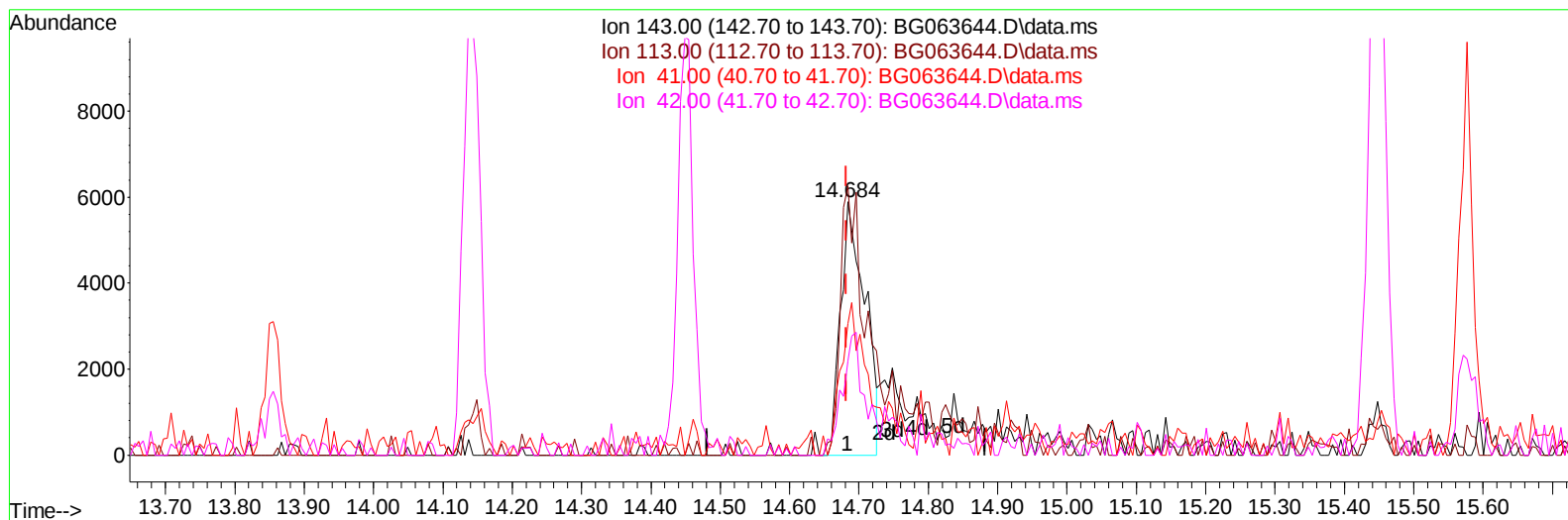


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
Data File : BG063644.D
Acq On : 10 Dec 2024 22:58
Operator : RC/JU
Sample : P5109-19
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Dec 10 23:54:19 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: BG063644.D\data.ms

(54) 4-Nitrophenol-d4 (S)

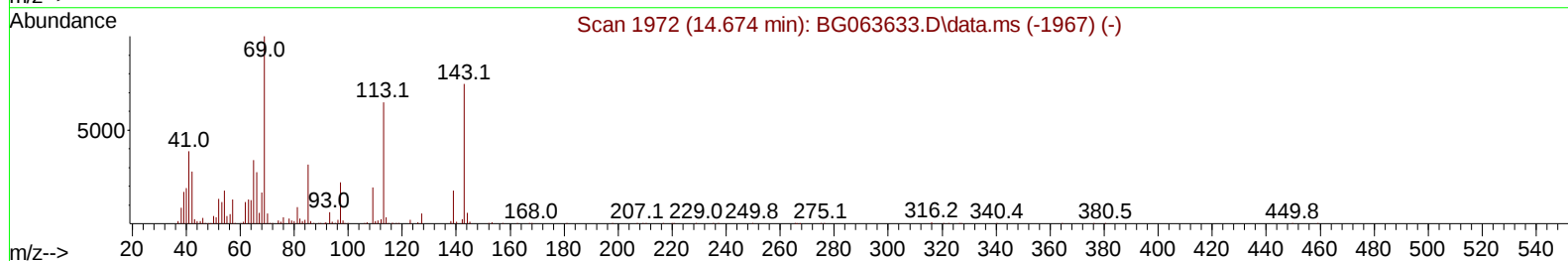
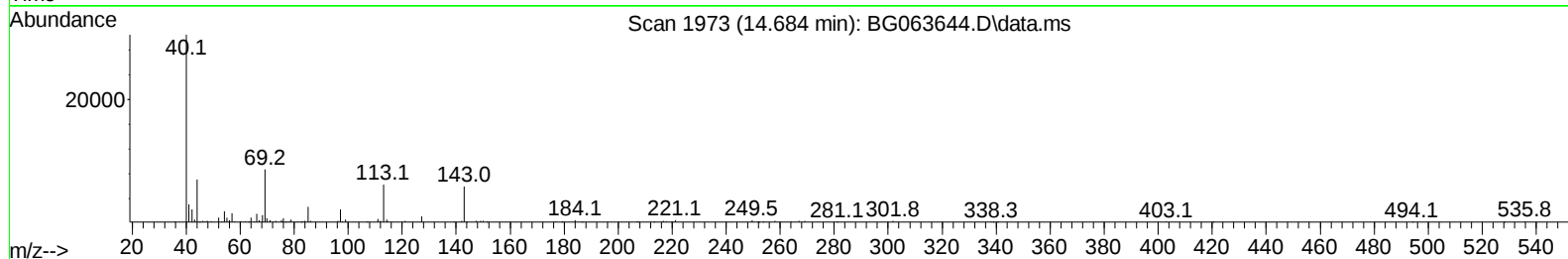
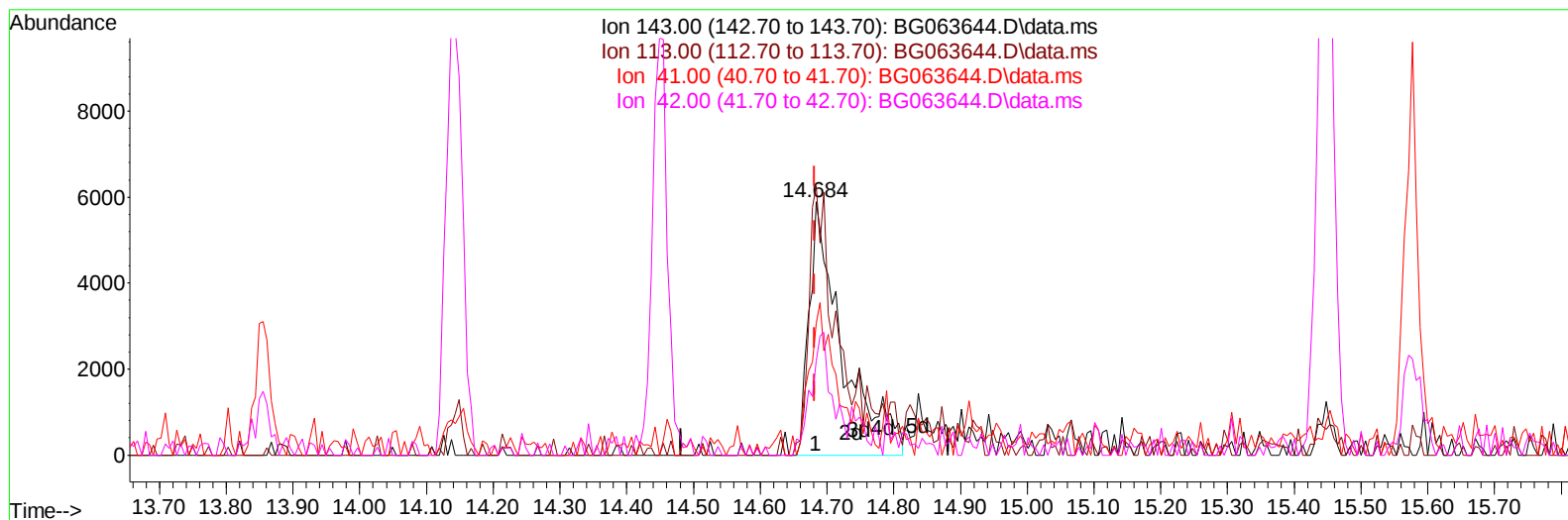
14.684min (+ 0.002) 3.62 ng/ul

response 14374

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	110.90	106.14
41.00	45.50	51.98
42.00	30.70	37.25#

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
Data File : BG063644.D
Acq On : 10 Dec 2024 22:58
Operator : RC/JU
Sample : P5109-19
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Dec 10 23:54:19 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: BG063644.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.684min (+ 0.002) 5.13 ng/ul m

response 20395

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	110.90	106.14
41.00	45.50	51.98
42.00	30.70	37.25#

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
 Data File : BG063644.D
 Acq On : 10 Dec 2024 22:58
 Operator : RC/JU
 Sample : P5109-19
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Dec 10 23:55:05 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	7.815	152	133462	20.000	ng/ul	0.00
20) Naphthalene-d8	10.600	136	563780	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.449	164	426908	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	1011396	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.452	240	979691	20.000	ng/ul	#-0.02
88) Perylene-d12	24.472	264	1078609	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.285	96	17062	5.644	ng/uL	0.00
4) Pyridine-d5	3.691	84	92490	8.270	ng/ul	0.00
7) Phenol-d5	6.993	99	93799	7.138	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.146	67	313552	35.291	ng/ul	0.00
11) 2-Chlorophenol-d4	7.351	132	248988	32.166	ng/ul	0.00
15) 4-Methylphenol-d8	8.526	113	187641	19.260	ng/ul	0.00
21) Nitrobenzene-d5	8.967	128	158672	40.096	ng/ul	0.00
24) 2-Nitrophenol-d4	9.684	143	171107	37.785	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.230	165	315029	32.309	ng/ul	0.00
31) 4-Chloroaniline-d4	10.741	131	332072	28.181	ng/ul	0.00
46) Dimethylphthalate-d6	13.855	166	1217478	39.240	ng/ul	0.00
49) Acenaphthylene-d8	14.143	160	1295353	38.195	ng/ul	0.00
54) 4-Nitrophenol-d4	14.684	143	20395m	5.134	ng/ul	0.00
60) Fluorene-d10	15.448	176	1082538	35.396	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.577	200	160638	28.115	ng/ul	0.00
73) Anthracene-d10	17.304	188	2003624	43.386	ng/ul	0.00
81) Pyrene-d10	19.602	212	2417088	44.511	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.273	264	2367892	43.710	ng/ul	-0.02

Target Compounds	Qvalue

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

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