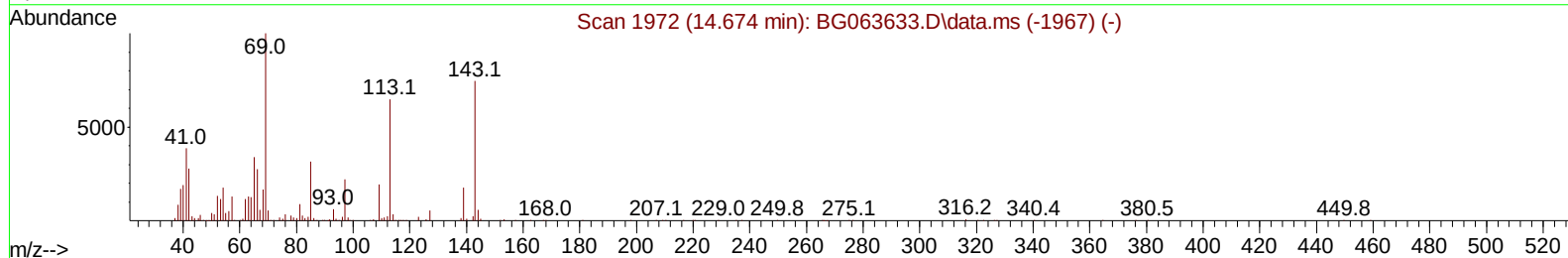
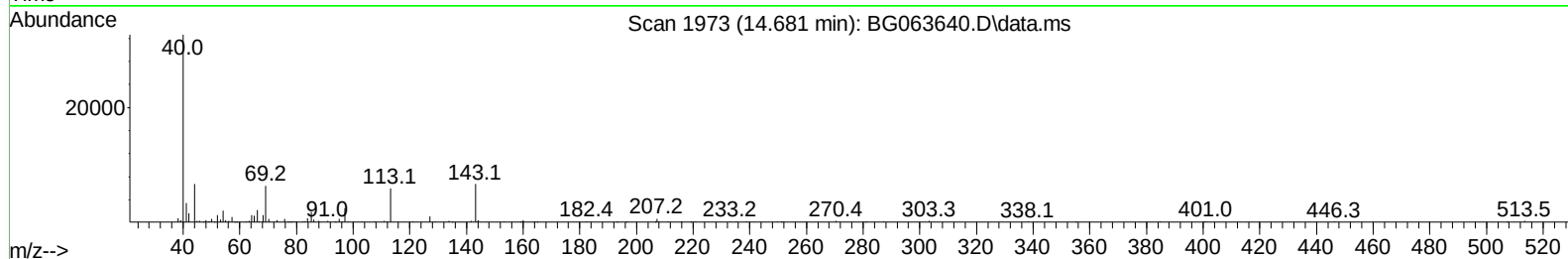
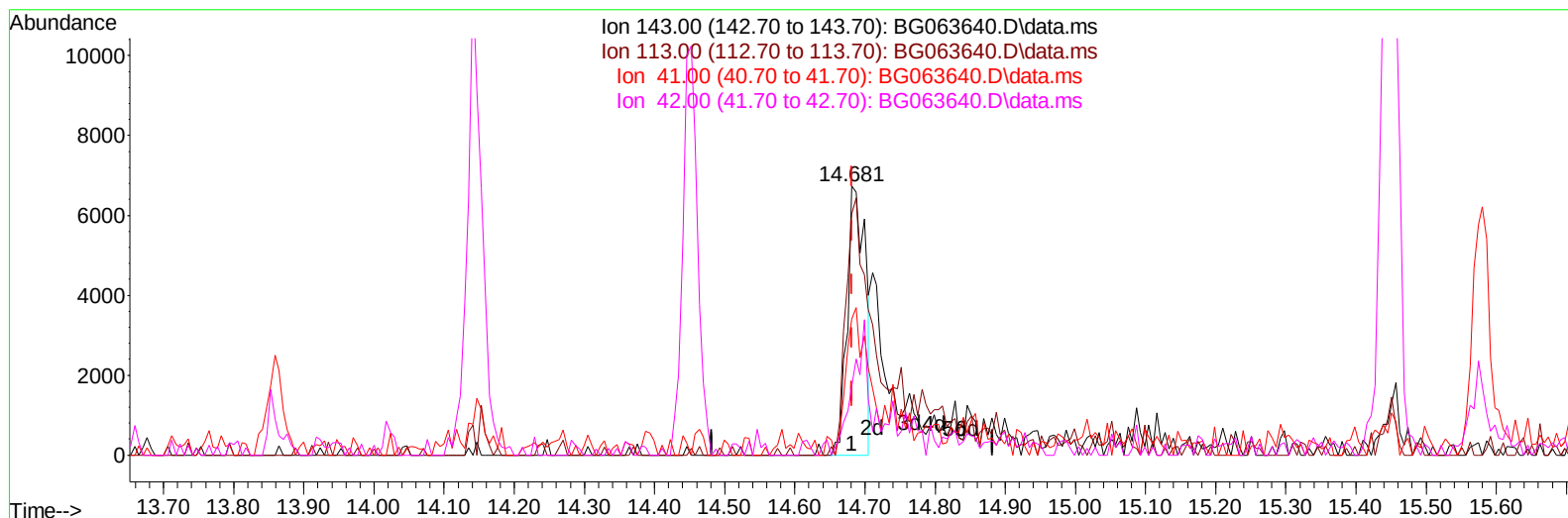


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
Data File : BG063640.D
Acq On : 10 Dec 2024 20:27
Operator : RC/JU
Sample : P5109-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Dec 10 23:27:10 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: BG063640.D\data.ms

(54) 4-Nitrophenol-d4 (S)

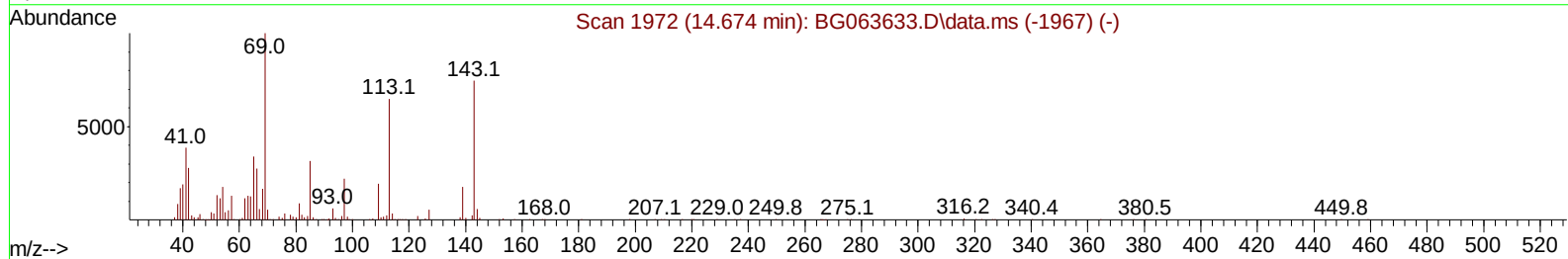
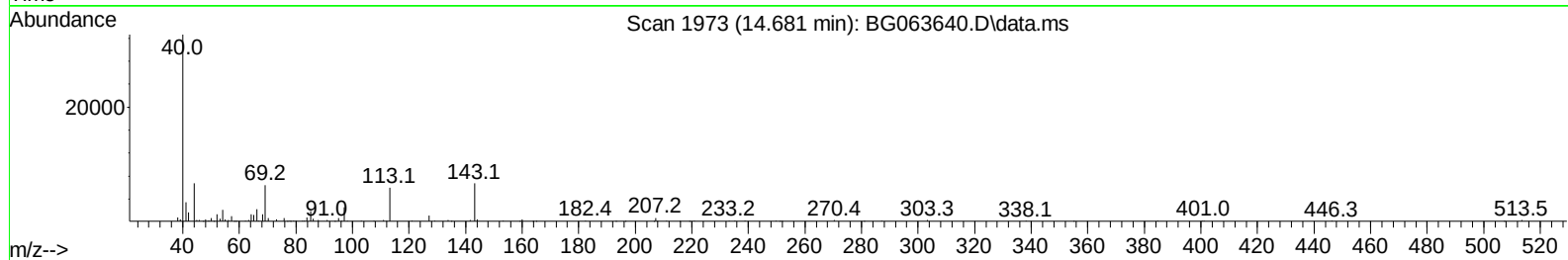
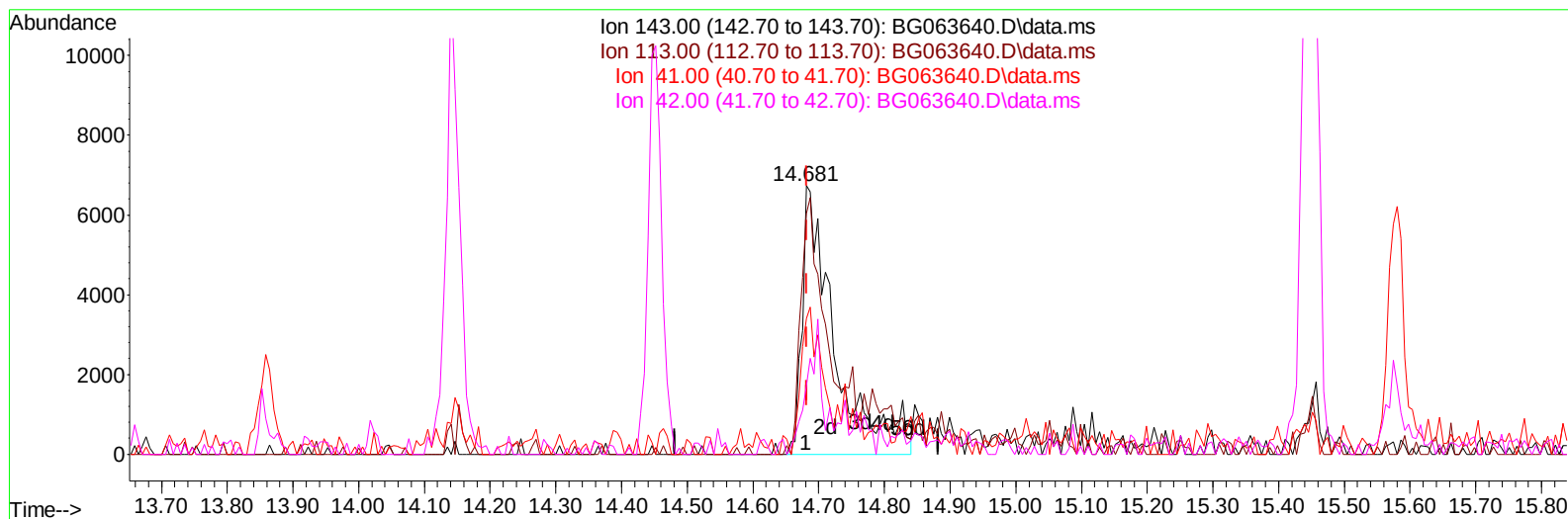
14.681min (-0.001) 2.89 ng/ul

response 12129

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	110.90	89.69
41.00	45.50	50.53
42.00	30.70	25.56

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
Data File : BG063640.D
Acq On : 10 Dec 2024 20:27
Operator : RC/JU
Sample : P5109-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Dec 10 23:27:10 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: BG063640.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.681min (-0.001) 5.48 ng/ul m

response 22986

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	110.90	89.69
41.00	45.50	50.53
42.00	30.70	25.56

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
 Data File : BG063640.D
 Acq On : 10 Dec 2024 20:27
 Operator : RC/JU
 Sample : P5109-15
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Dec 10 23:44:26 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	7.813	152	140716	20.000	ng/ul	0.00
20) Naphthalene-d8	10.598	136	618332	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.452	164	450719	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.202	188	1053715	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.456	240	995041	20.000	ng/ul	#-0.02
88) Perylene-d12	24.476	264	1074389	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.283	96	17483	5.485	ng/uL	0.00
4) Pyridine-d5	3.688	84	99205	8.413	ng/ul	0.00
7) Phenol-d5	6.996	99	104138	7.517	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.143	67	301090	32.142	ng/ul	0.00
11) 2-Chlorophenol-d4	7.349	132	249138	30.526	ng/ul	0.00
15) 4-Methylphenol-d8	8.530	113	197914	19.267	ng/ul	0.00
21) Nitrobenzene-d5	8.970	128	144526	33.299	ng/ul	0.00
24) 2-Nitrophenol-d4	9.687	143	164239	33.068	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	306729	28.683	ng/ul	0.00
31) 4-Chloroaniline-d4	10.739	131	374977	29.015	ng/ul	0.00
46) Dimethylphthalate-d6	13.859	166	1121197	34.228	ng/ul	0.00
49) Acenaphthylene-d8	14.141	160	1188131	33.183	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.681	143	22986m	5.480	ng/ul	0.00
60) Fluorene-d10	15.451	176	1001296	31.010	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.580	200	129660	21.782	ng/ul	0.00
73) Anthracene-d10	17.302	188	1861881	38.697	ng/ul	-0.01
81) Pyrene-d10	19.605	212	2125142	38.531	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.270	264	2122848	39.340	ng/ul	-0.02

Target Compounds	Qvalue

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

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