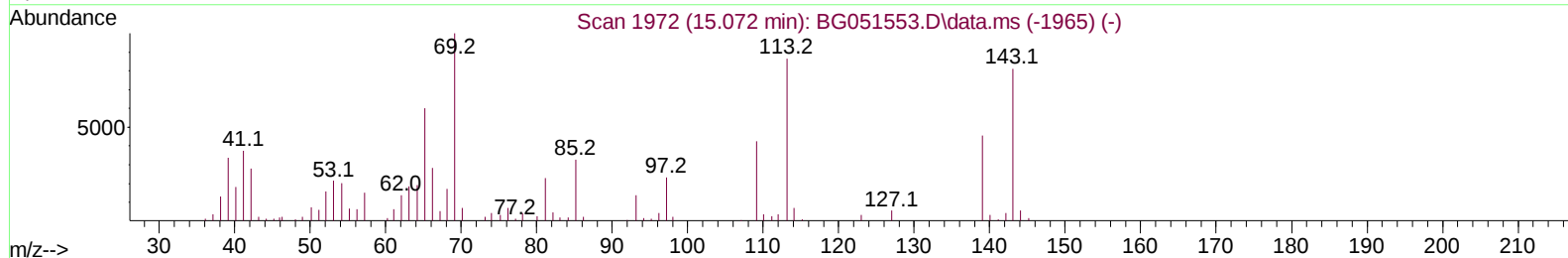
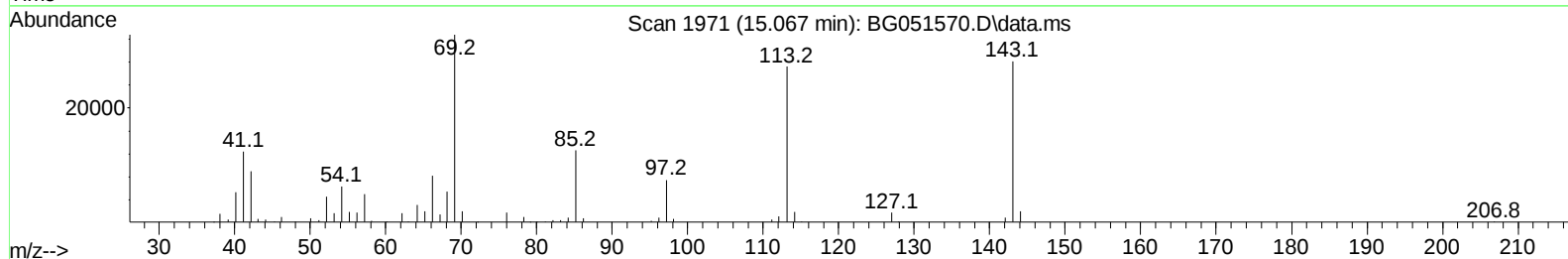
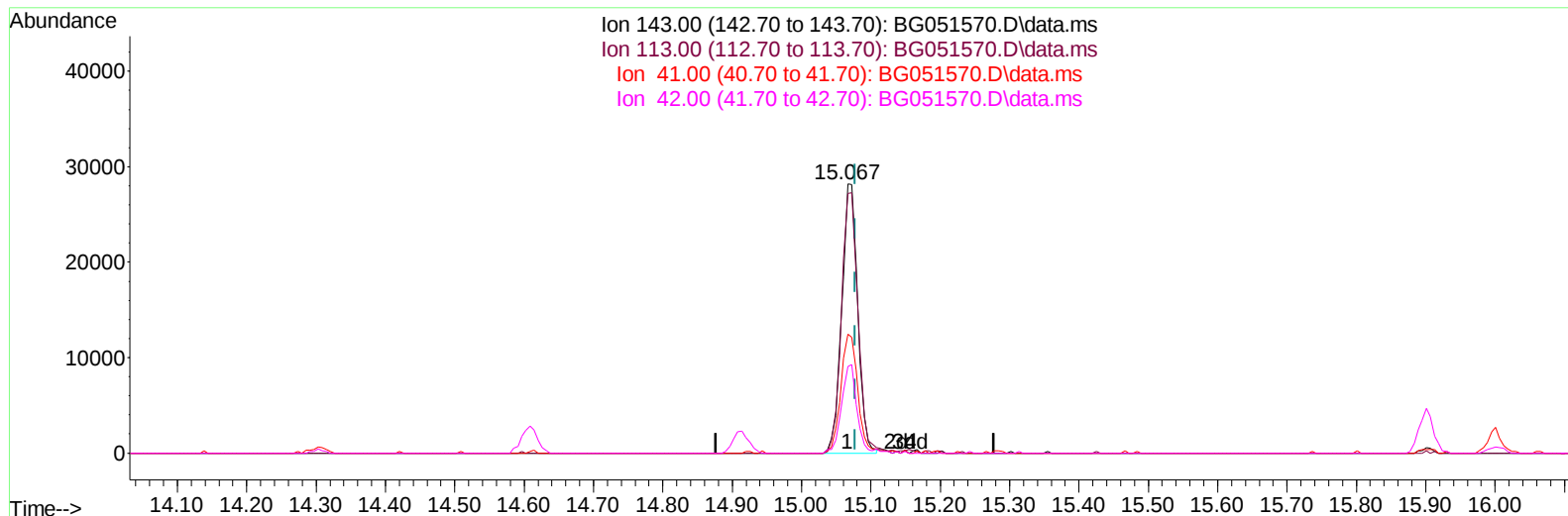


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051570.D
Acq On : 16 Dec 2021 14:46
Operator : CG/JU
Sample : PB141433BL
Misc :
ALS Vial : 70 Sample Multiplier: 1

Quant Time: Dec 17 00:11:21 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 14 14:19:57 2021
Response via : Initial Calibration



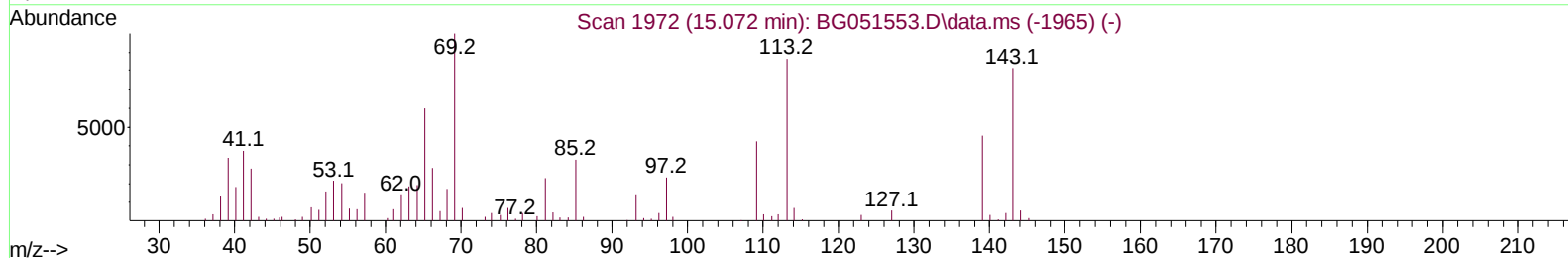
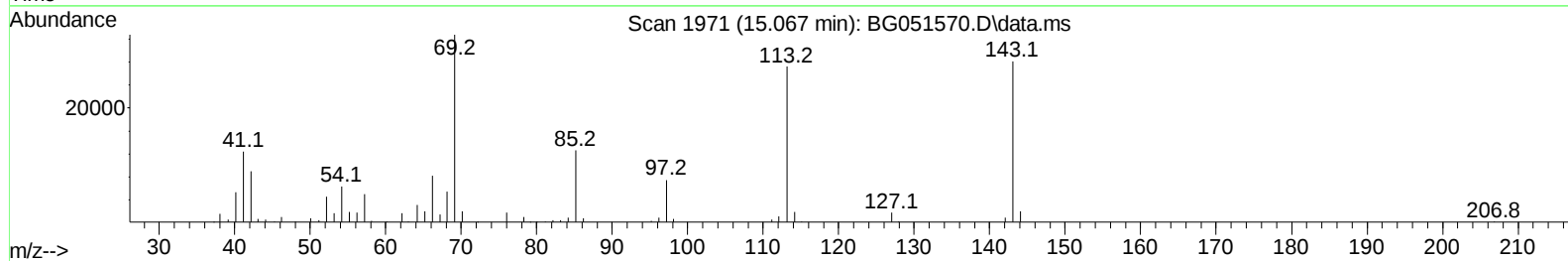
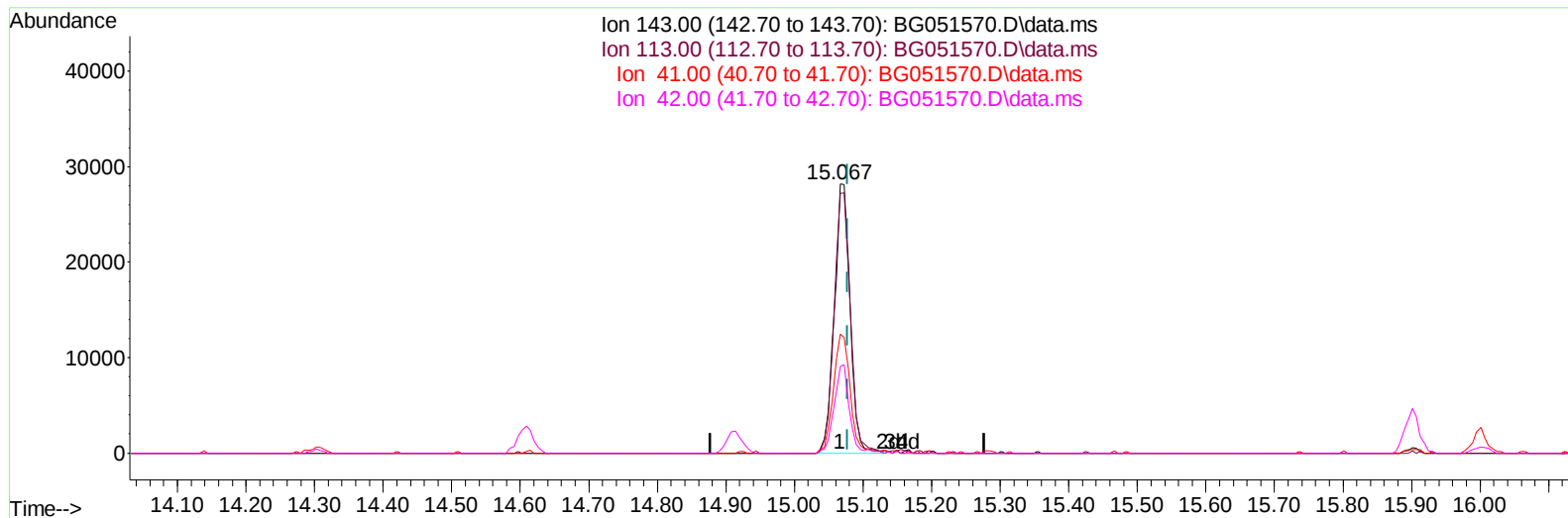
TIC: BG051570.D\data.ms

(54) 4-Nitrophenol-d4 (S)**15.067min (-0.011) 33.14 ng/u1****response 44812**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	96.54#
41.00	44.40	44.21
42.00	29.70	32.11

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
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Acq On : 16 Dec 2021 14:46
Operator : CG/JU
Sample : PB141433BL
Misc :
ALS Vial : 70 Sample Multiplier: 1

Quant Time: Dec 17 00:11:21 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
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Response via : Initial Calibration



TIC: BG051570.D\data.ms

(54) 4-Nitrophenol-d4 (S)**15.067min (-0.011) 33.38 ng/ul m****response 45126**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	96.54#
41.00	44.40	44.21
42.00	29.70	32.11

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
 Data File : BG051570.D
 Acq On : 16 Dec 2021 14:46
 Operator : CG/JU
 Sample : PB141433BL
 Misc :
 ALS Vial : 70 Sample Multiplier: 1

Quant Time: Dec 17 00:11:21 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.276	152	33843	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.119	136	137695	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.914	164	103877	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.657	188	256639	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.981	240	260861	20.000	ng/ul	# 0.00
88) Perylene-d12	25.476	264	255403	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.606	96	5157	6.328	ng/uL	0.00
4) Pyridine-d5	4.029	84	79908	32.454	ng/ul	-0.02
7) Phenol-d5	7.406	99	101292	33.521	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.583	67	59908	33.350	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.800	132	71993	34.256	ng/ul	-0.01
15) 4-Methylphenol-d8	8.975	113	79057	33.847	ng/ul	0.00
21) Nitrobenzene-d5	9.445	128	36741	36.394	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.179	143	41839	37.912	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.720	165	81347	33.991	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.237	131	106687	33.859	ng/ul	0.00
46) Dimethylphthalate-d6	14.303	166	290454	34.934	ng/ul	-0.01
49) Acenaphthylene-d8	14.608	160	347925	33.694	ng/ul	0.00
54) 4-Nitrophenol-d4	15.067	143	45126m	33.376	ng/ul	-0.01
60) Fluorene-d10	15.901	176	247788	33.265	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.001	200	43838	32.786	ng/ul	-0.01
73) Anthracene-d10	17.757	188	423231	34.571	ng/ul	0.00
81) Pyrene-d10	20.031	212	526042	33.506	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.241	264	472984	35.197	ng/ul	-0.01

Target Compounds	Qvalue

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

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