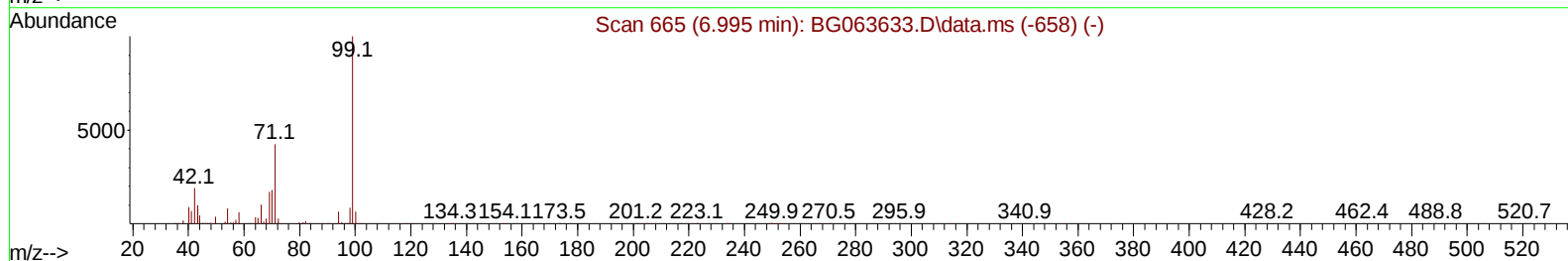
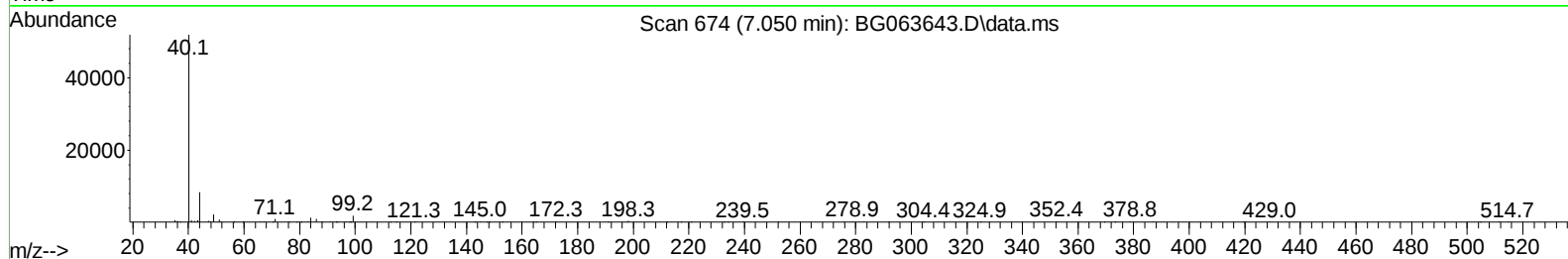
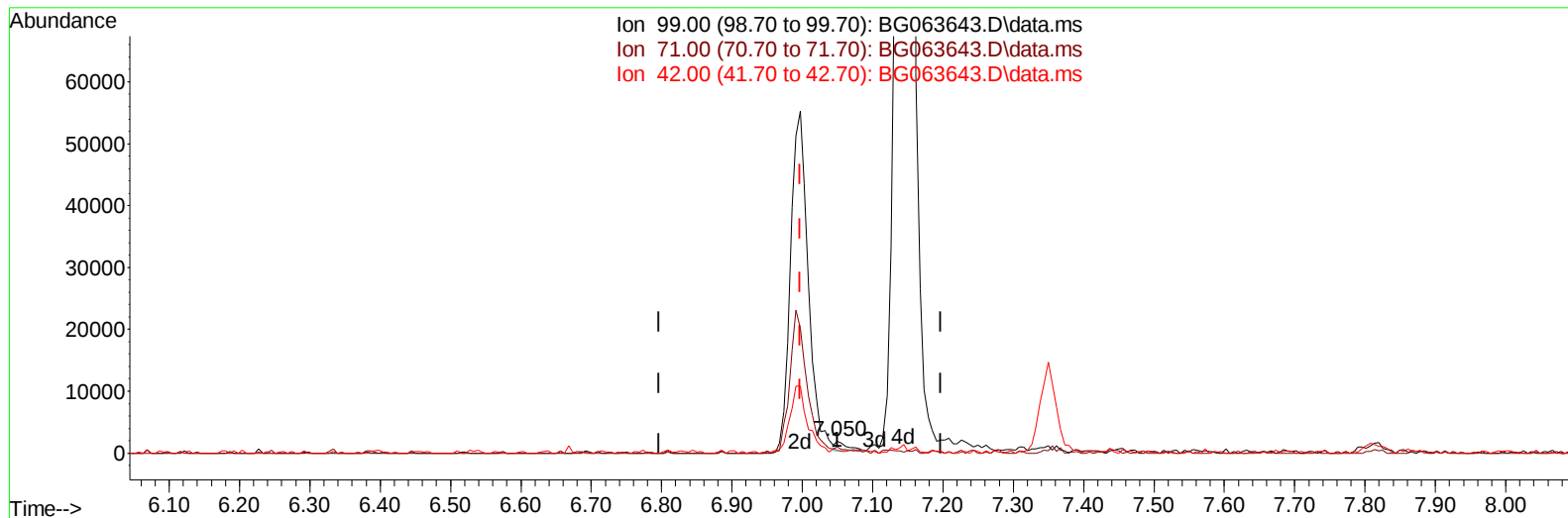


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
Data File : BG063643.D
Acq On : 10 Dec 2024 22:20
Operator : RC/JU
Sample : P5109-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Quant Time: Dec 10 23:27:37 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: BG063643.D\data.ms

(7) Phenol-d5 (S)

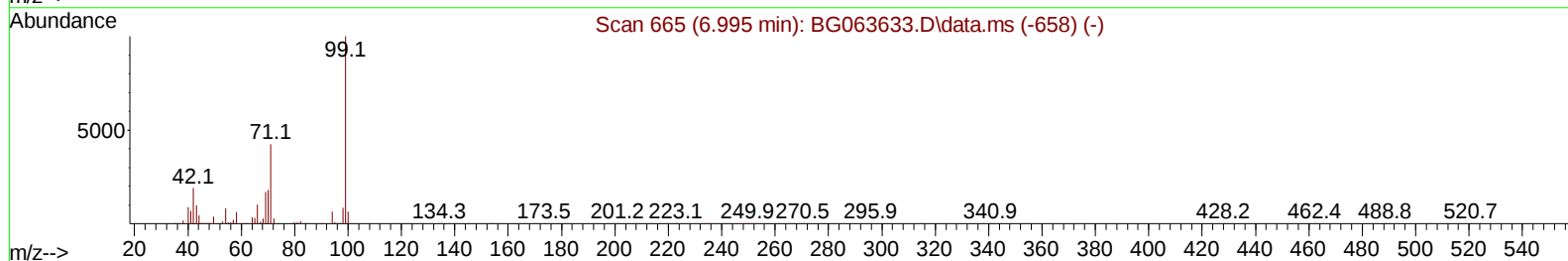
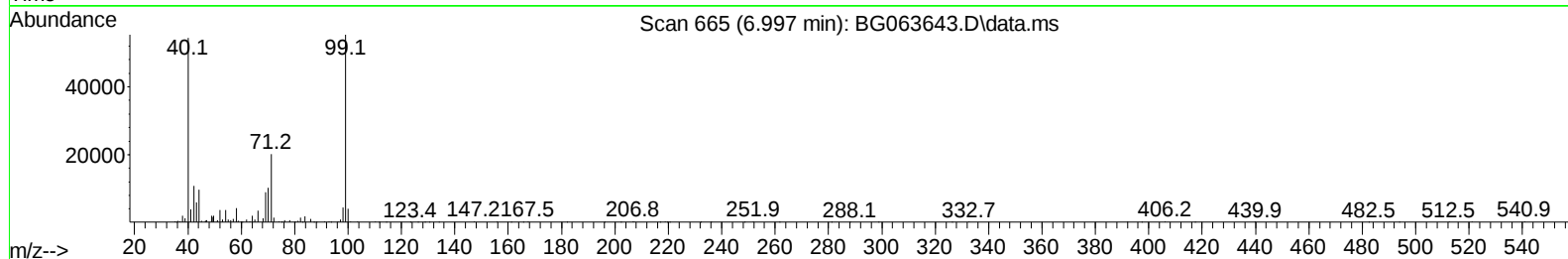
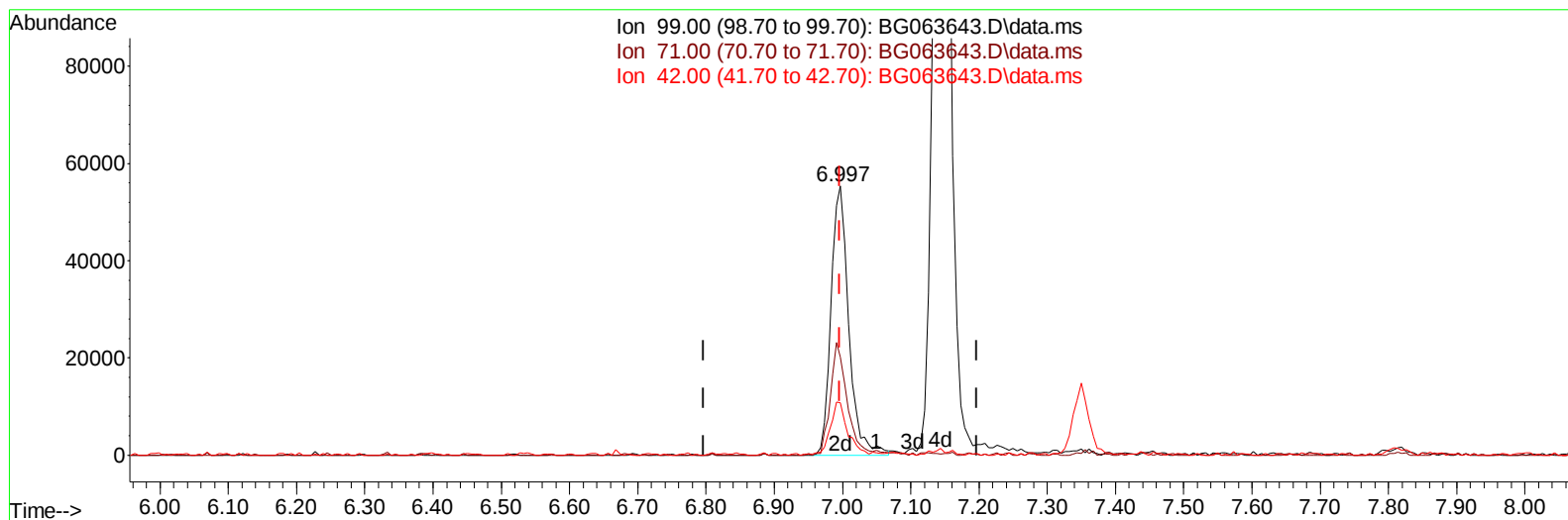
7.050min (+ 0.053) 0.13 ng/u1

response 1888

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	54.50	50.69
42.00	24.30	22.05
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
Data File : BG063643.D
Acq On : 10 Dec 2024 22:20
Operator : RC/JU
Sample : P5109-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Quant Time: Dec 10 23:27:37 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: BG063643.D\data.ms

(7) Phenol-d5 (S)

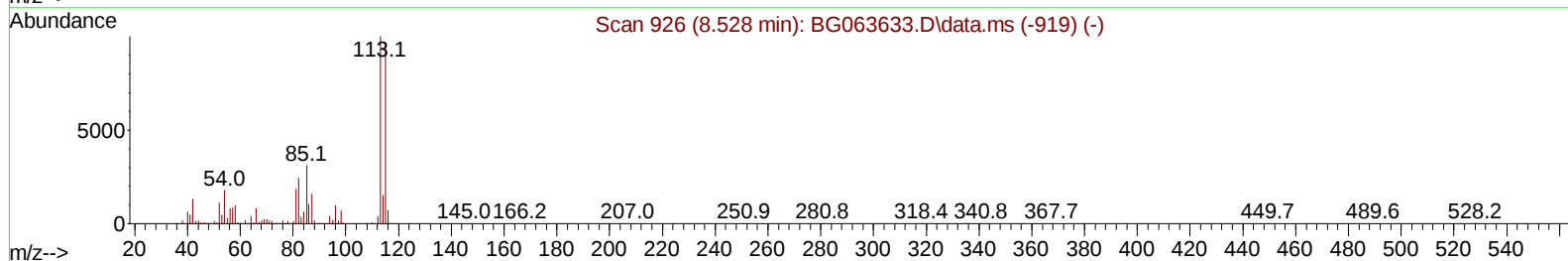
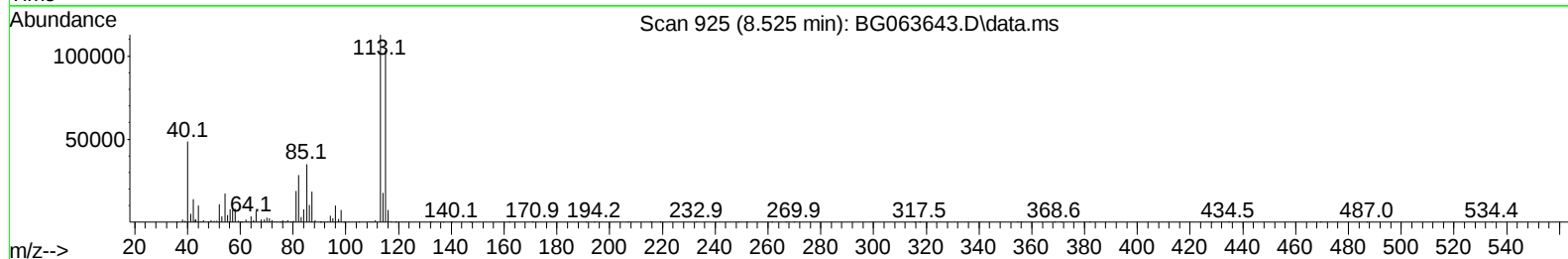
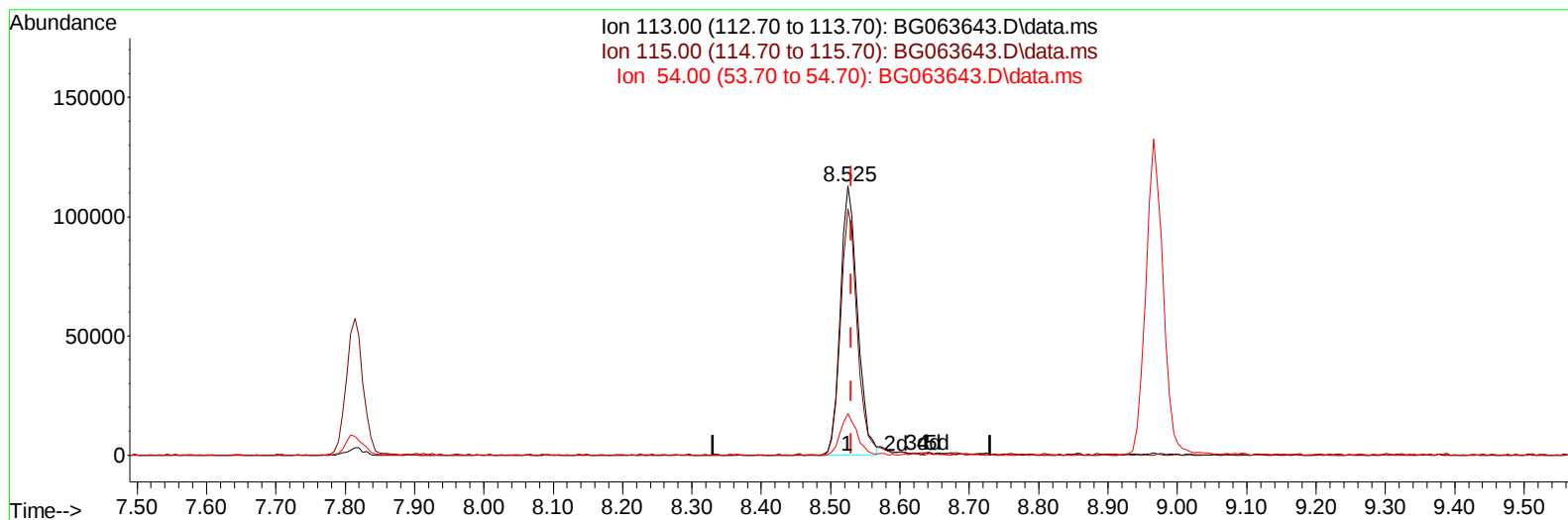
6.997min (+ 0.000) 6.60 ng/u1 m

response 99496

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	54.50	36.45#
42.00	24.30	19.58
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
Data File : BG063643.D
Acq On : 10 Dec 2024 22:20
Operator : RC/JU
Sample : P5109-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Quant Time: Dec 10 23:27:37 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
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TIC: BG063643.D\data.ms

(15) 4-Methylphenol-d8 (S)

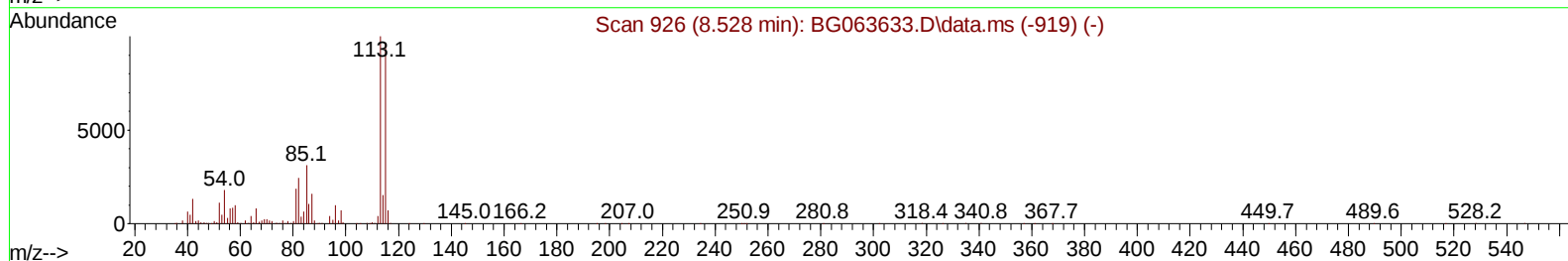
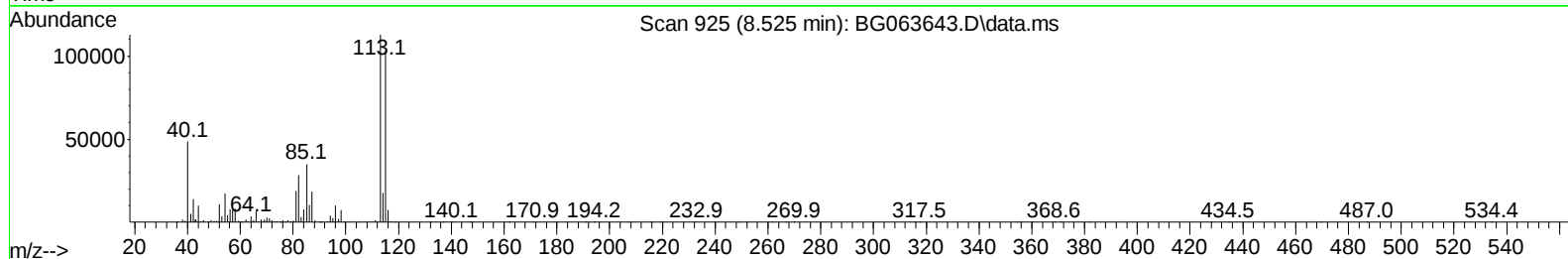
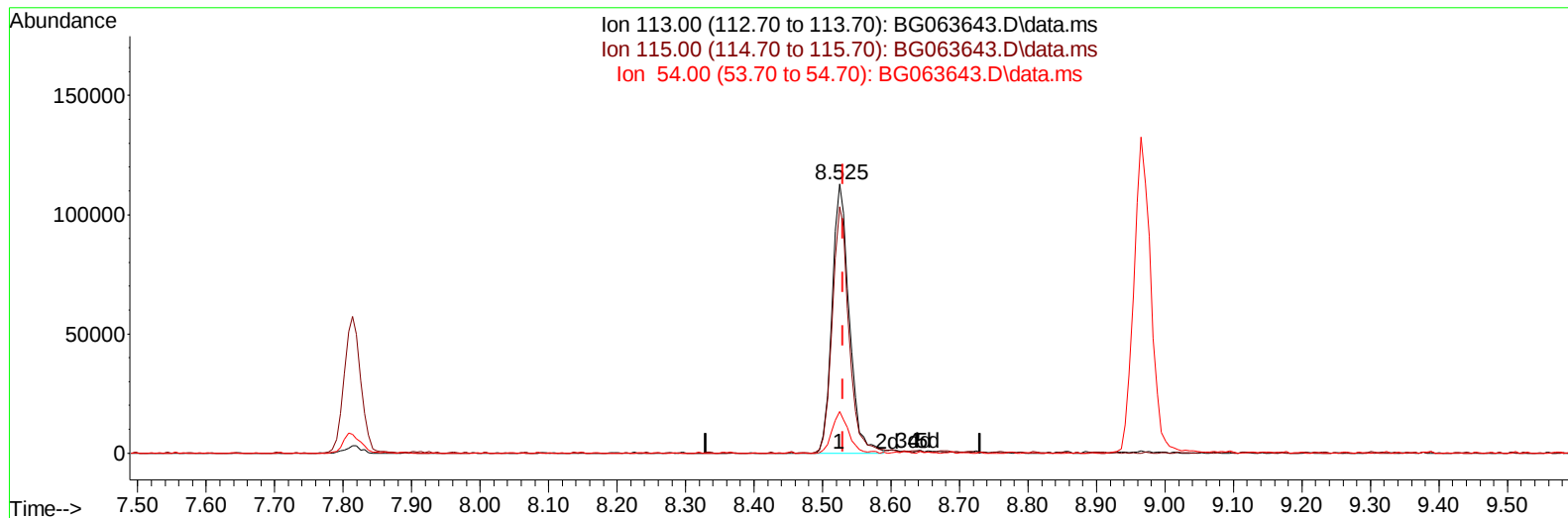
8.525min (-0.006) 17.29 ng/u1

response 193282

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	86.90	91.57
54.00	18.30	15.44
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
Data File : BG063643.D
Acq On : 10 Dec 2024 22:20
Operator : RC/JU
Sample : P5109-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Quant Time: Dec 10 23:27:37 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: BG063643.D\data.ms

(15) 4-Methylphenol-d8 (S)

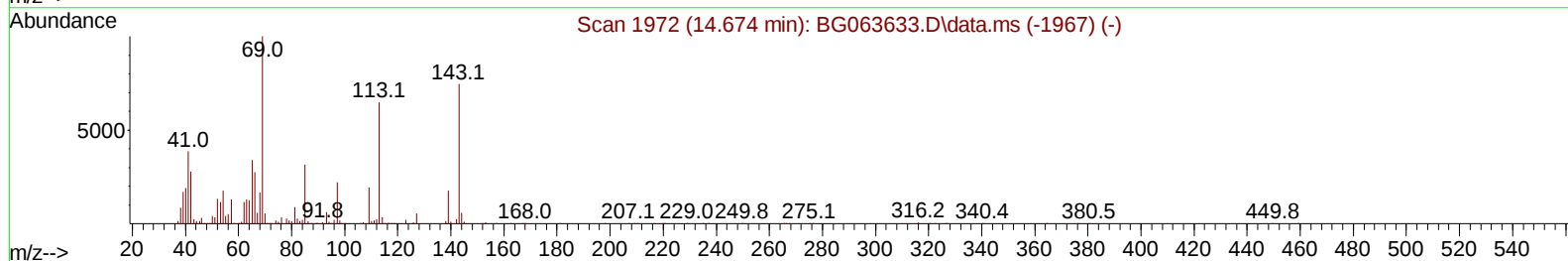
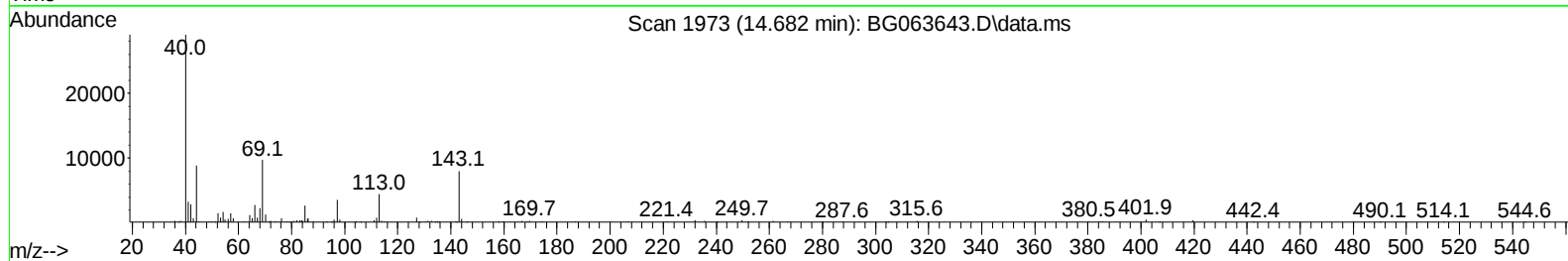
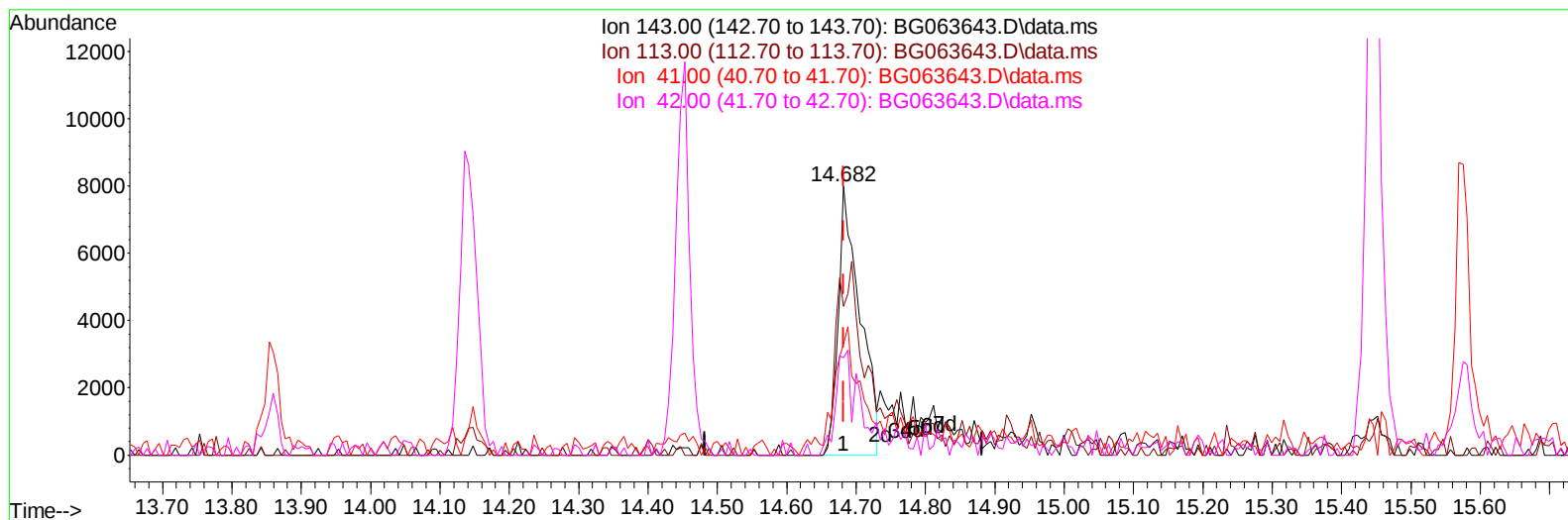
8.525min (-0.006) 17.57 ng/ul m

response 196448

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	86.90	91.57
54.00	18.30	15.44
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
Data File : BG063643.D
Acq On : 10 Dec 2024 22:20
Operator : RC/JU
Sample : P5109-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Quant Time: Dec 10 23:27:37 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120624.MA.M
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TIC: BG063643.D\data.ms

(54) 4-Nitrophenol-d4 (S)

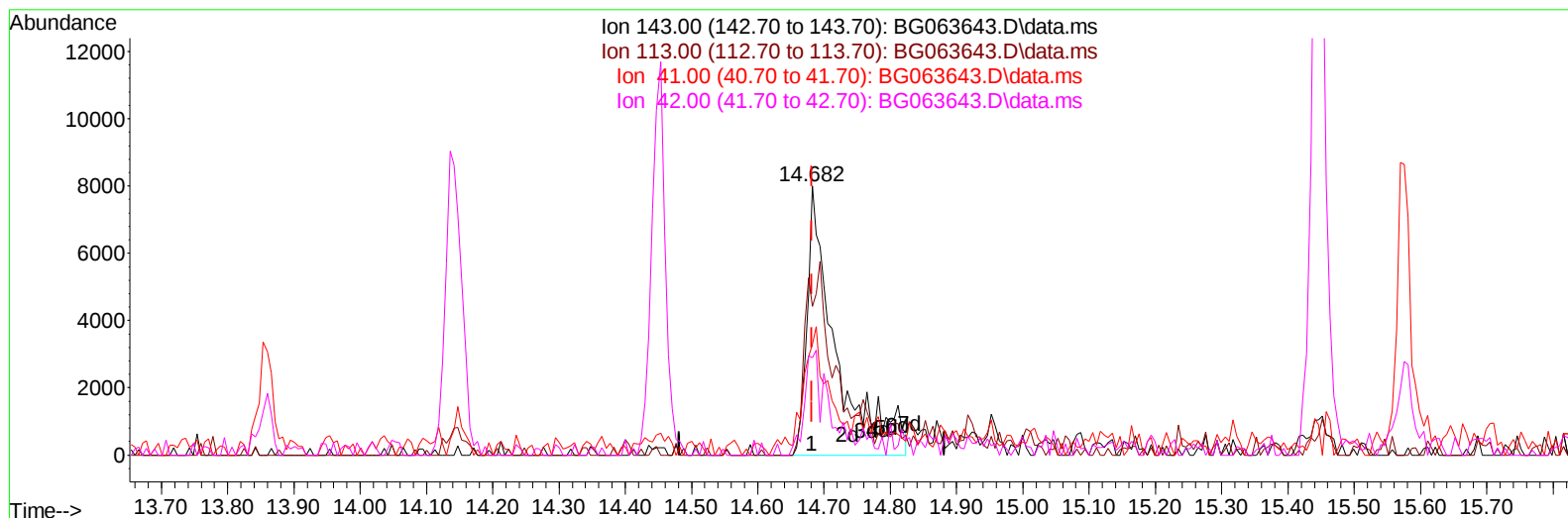
14.682min (+ 0.000) 3.83 ng/ul

response 17417

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	110.90	55.33#
41.00	45.50	41.27
42.00	30.70	36.09

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
Data File : BG063643.D
Acq On : 10 Dec 2024 22:20
Operator : RC/JU
Sample : P5109-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Quant Time: Dec 10 23:27:37 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120624.MA.M
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Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
 Data File : BG063643.D
 Acq On : 10 Dec 2024 22:20
 Operator : RC/JU
 Sample : P5109-18
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Dec 10 23:49:38 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	7.814	152	153126	20.000	ng/ul	0.00
20) Naphthalene-d8	10.599	136	652951	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.453	164	489192	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.203	188	1115418	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.457	240	1048621	20.000	ng/ul	#-0.02
88) Perylene-d12	24.477	264	1112498	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.278	96	18377	5.298	ng/uL	0.00
4) Pyridine-d5	3.689	84	90047	7.018	ng/ul	0.00
7) Phenol-d5	6.997	99	99496m	6.599	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.144	67	298286	29.262	ng/ul	0.00
11) 2-Chlorophenol-d4	7.350	132	245748	27.671	ng/ul	0.00
15) 4-Methylphenol-d8	8.525	113	196448m	17.574	ng/ul	0.00
21) Nitrobenzene-d5	8.971	128	145273	31.697	ng/ul	0.00
24) 2-Nitrophenol-d4	9.688	143	164077	31.284	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.229	165	316256	28.006	ng/ul	0.00
31) 4-Chloroaniline-d4	10.740	131	316210	23.171	ng/ul	0.00
46) Dimethylphthalate-d6	13.860	166	1167349	32.834	ng/ul	0.00
49) Acenaphthylene-d8	14.142	160	1205756	31.027	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.682	143	24225m	5.321	ng/ul	0.00
60) Fluorene-d10	15.446	176	1011622	28.866	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.581	200	146114	23.188	ng/ul	0.00
73) Anthracene-d10	17.303	188	1927152	37.838	ng/ul	-0.01
81) Pyrene-d10	19.600	212	2232669	38.412	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.271	264	2188654	39.170	ng/ul	-0.02

Target Compounds	Qvalue

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

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