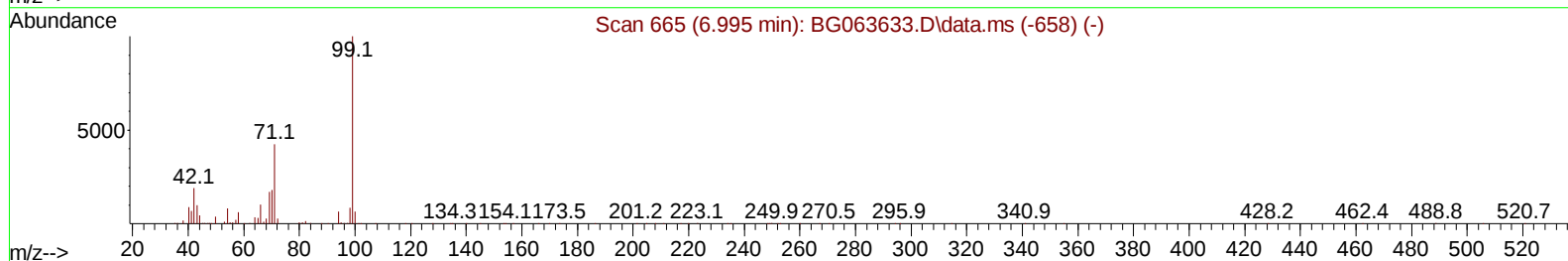
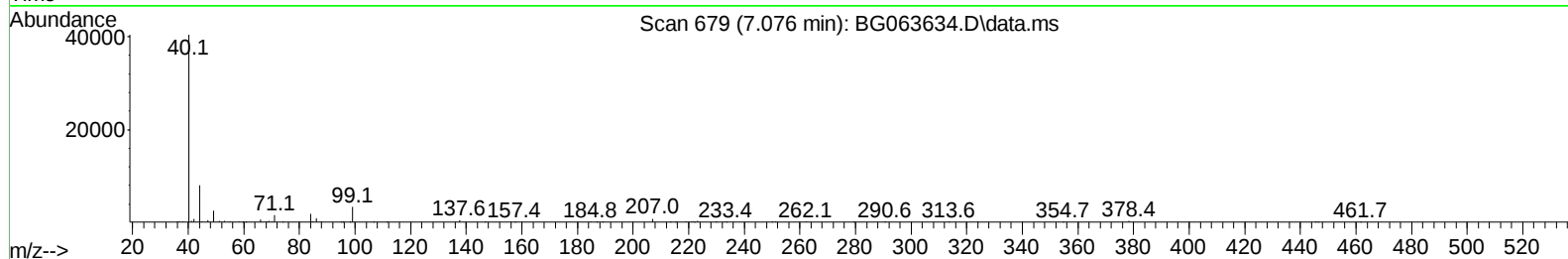
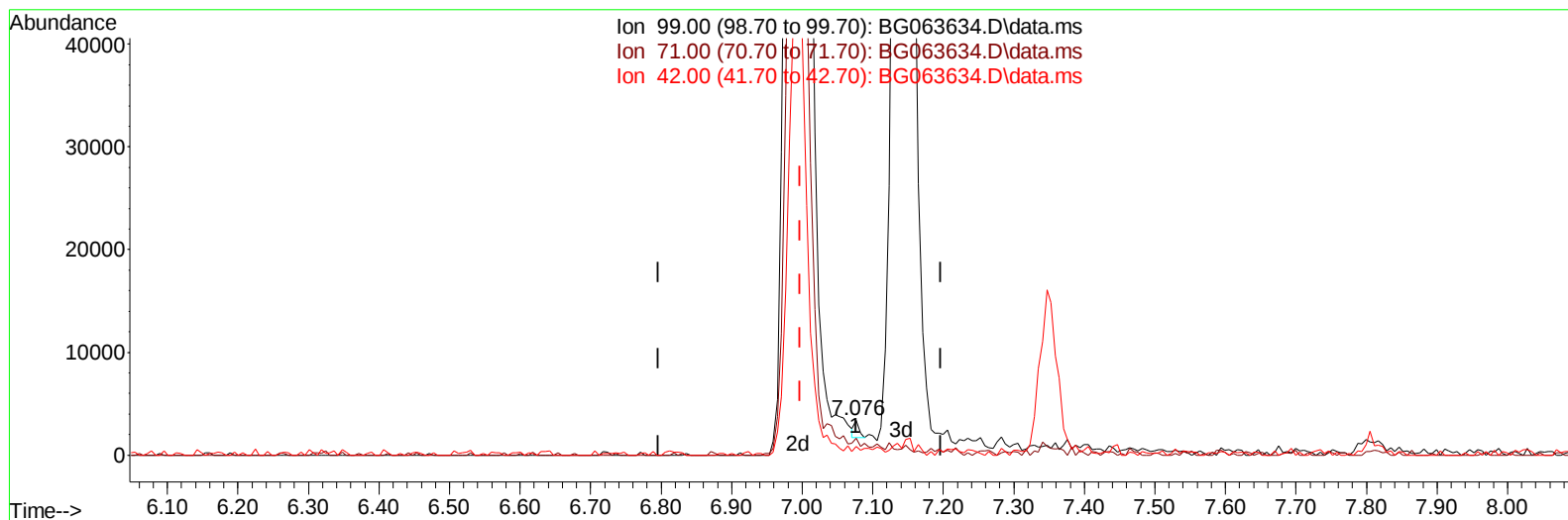


Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
Data File : BG063634.D
Acq On : 10 Dec 2024 16:41
Operator : RC/JU
Sample : PB165508BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Dec 10 17:32:31 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: BG063634.D\data.ms

(7) Phenol-d5 (S)

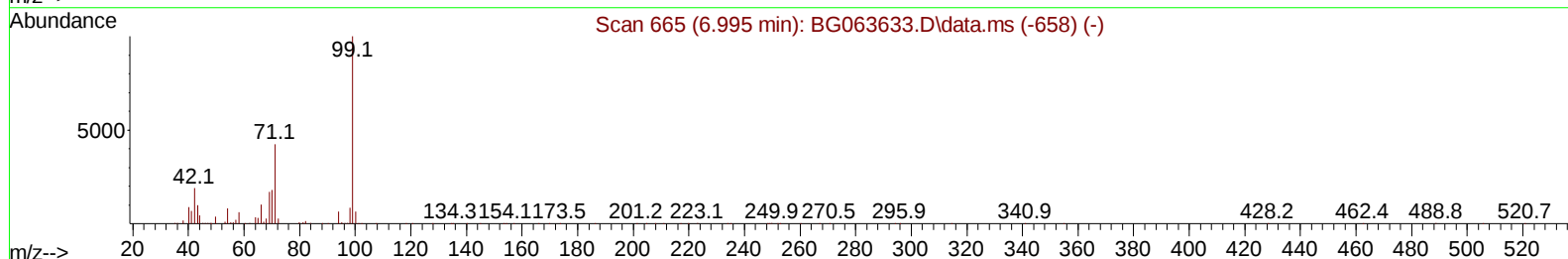
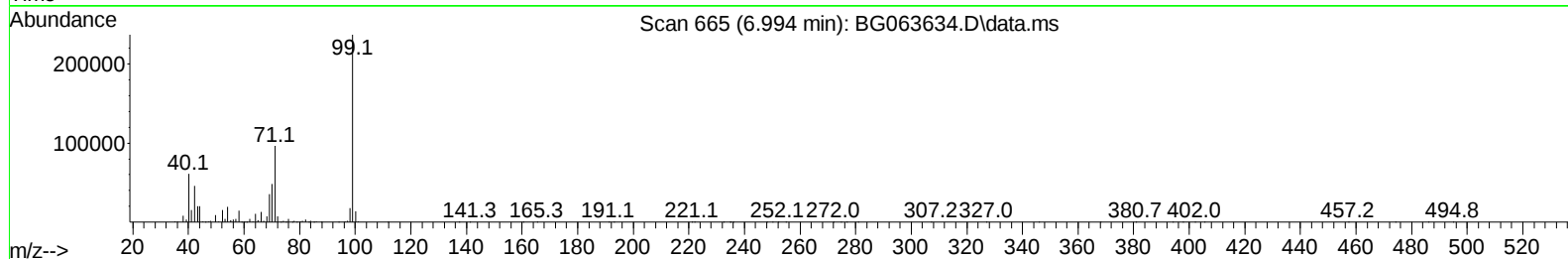
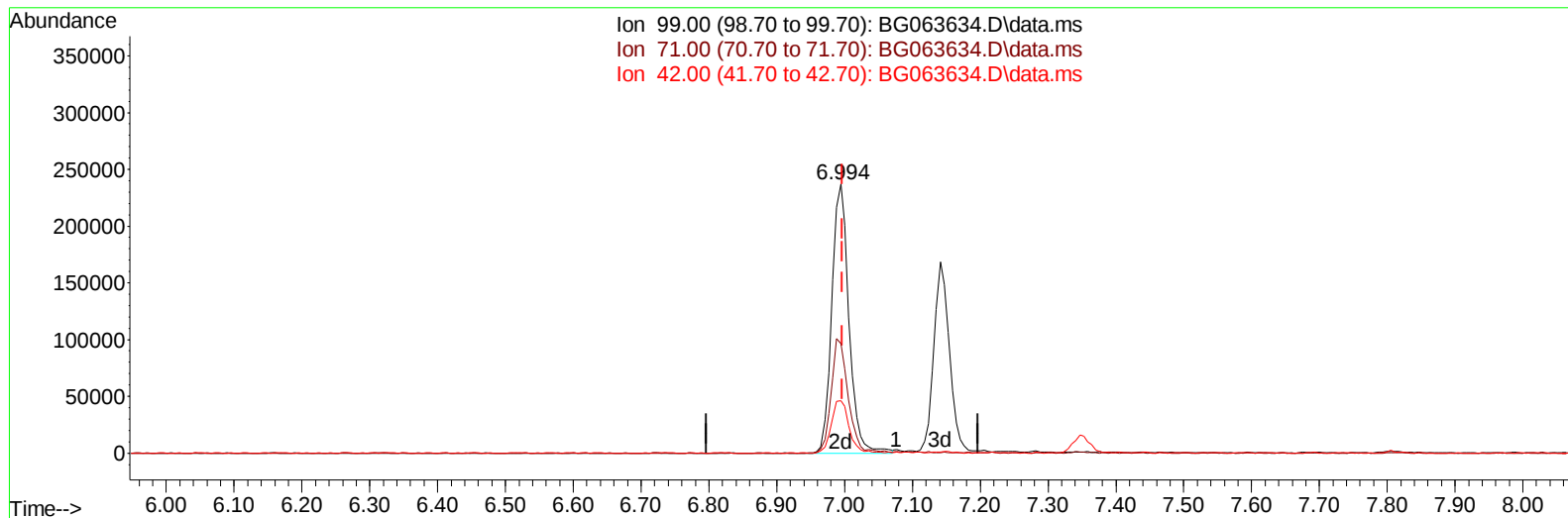
7.076min (+ 0.080) 0.06 ng/u1

response 740

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	54.50	48.34
42.00	24.30	24.24
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
Data File : BG063634.D
Acq On : 10 Dec 2024 16:41
Operator : RC/JU
Sample : PB165508BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Dec 10 17:32:31 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: BG063634.D\data.ms

(7) Phenol-d5 (S)

6.994min (-0.003) 31.47 ng/ul m

response 416679

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	54.50	40.62#
42.00	24.30	19.48
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
 Data File : BG063634.D
 Acq On : 10 Dec 2024 16:41
 Operator : RC/JU
 Sample : PB165508BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Dec 10 17:33:00 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	7.811	152	134476	20.000	ng/ul	0.00
20) Naphthalene-d8	10.602	136	579083	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.456	164	439736	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.206	188	1032679	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.460	240	929349	20.000	ng/ul	#-0.01
88) Perylene-d12	24.485	264	1007738	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	21861	7.177	ng/uL	0.00
4) Pyridine-d5	3.686	84	325289	28.867	ng/ul	0.00
7) Phenol-d5	6.994	99	416679m	31.471	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.141	67	262151	29.283	ng/ul	0.00
11) 2-Chlorophenol-d4	7.347	132	278609	35.722	ng/ul	0.00
15) 4-Methylphenol-d8	8.528	113	329957	33.612	ng/ul	0.00
21) Nitrobenzene-d5	8.968	128	134899	33.188	ng/ul	0.00
24) 2-Nitrophenol-d4	9.685	143	147379	31.685	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.232	165	306947	30.648	ng/ul	0.00
31) 4-Chloroaniline-d4	10.737	131	375390	31.016	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	994057	31.104	ng/ul	0.00
49) Acenaphthylene-d8	14.145	160	1169383	33.475	ng/ul	0.00
54) 4-Nitrophenol-d4	14.679	143	102494	25.046	ng/ul	0.00
60) Fluorene-d10	15.449	176	906537	28.776	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.584	200	84788	14.534	ng/ul	0.00
73) Anthracene-d10	17.306	188	1541854	32.699	ng/ul	0.00
81) Pyrene-d10	19.603	212	1768031	34.322	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.274	264	1649429	32.589	ng/ul	-0.02

Target Compounds	Qvalue

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

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