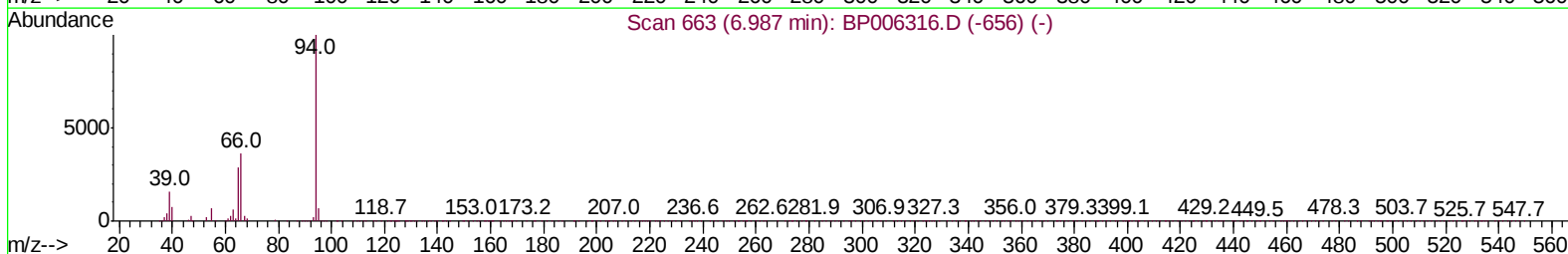
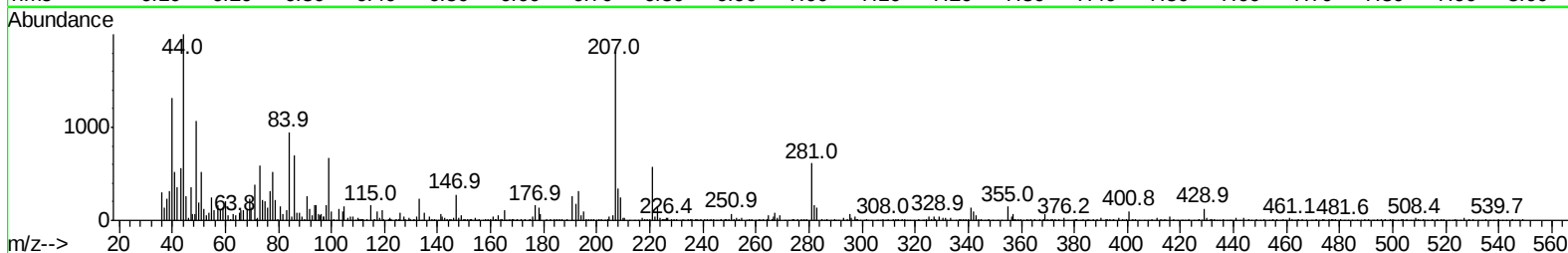
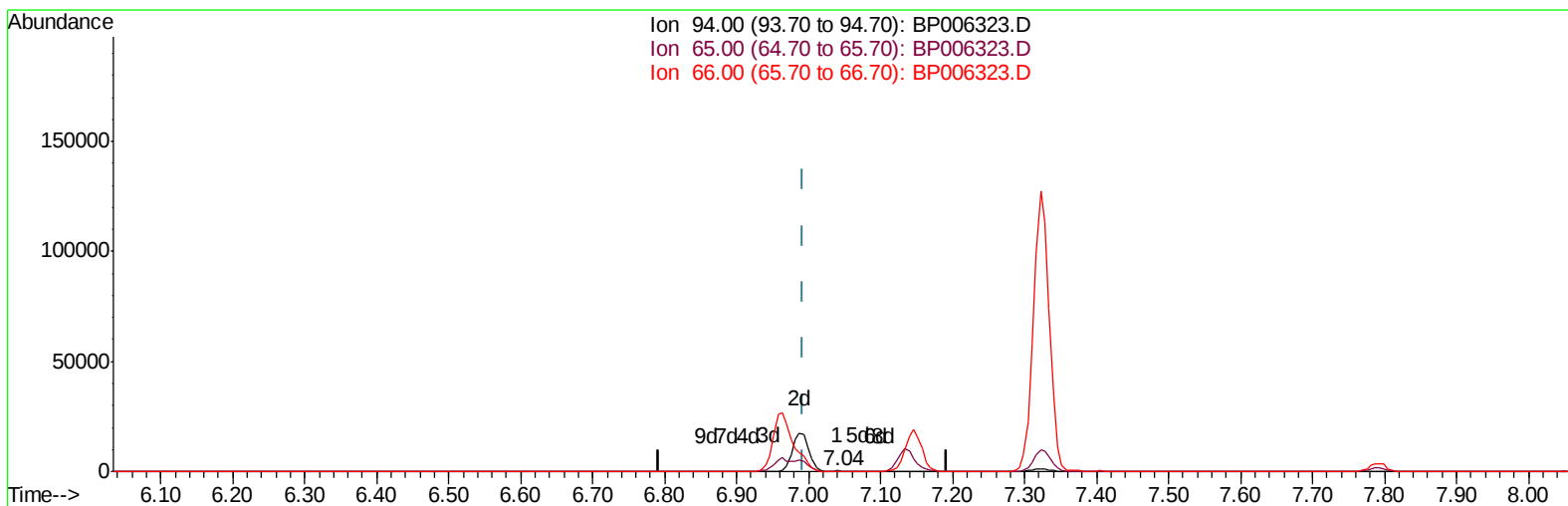


Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP072421\
Data File : BP006323.D
Acq On : 25 Jul 2021 01:21
Operator : CG/JU
Sample : M2973-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Quant Time: Jul 26 01:10:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jul 24 04:29:12 2021
Response via : Initial Calibration



TIC: BP006323.D

(8) Phenol

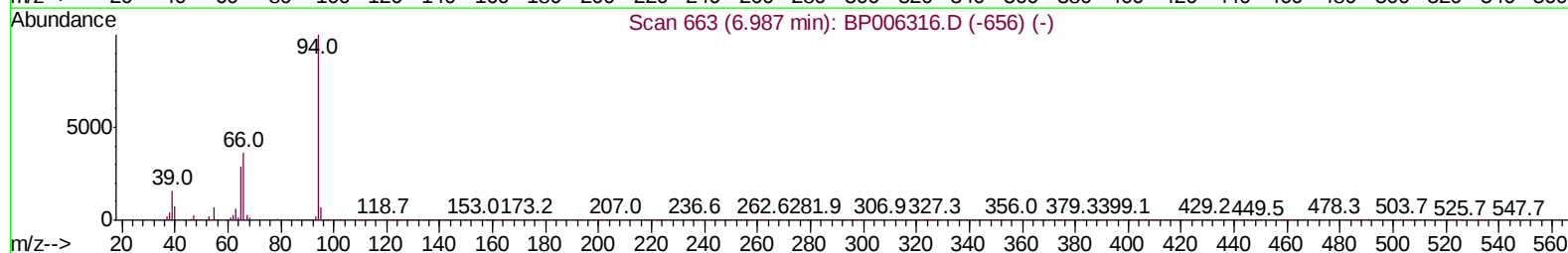
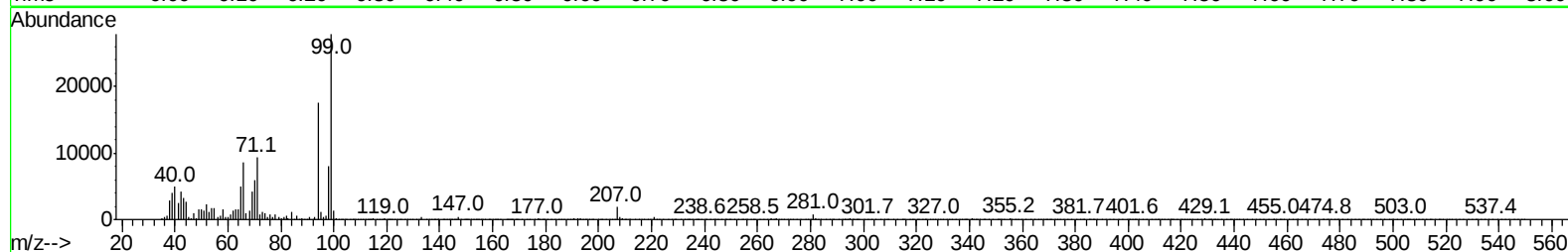
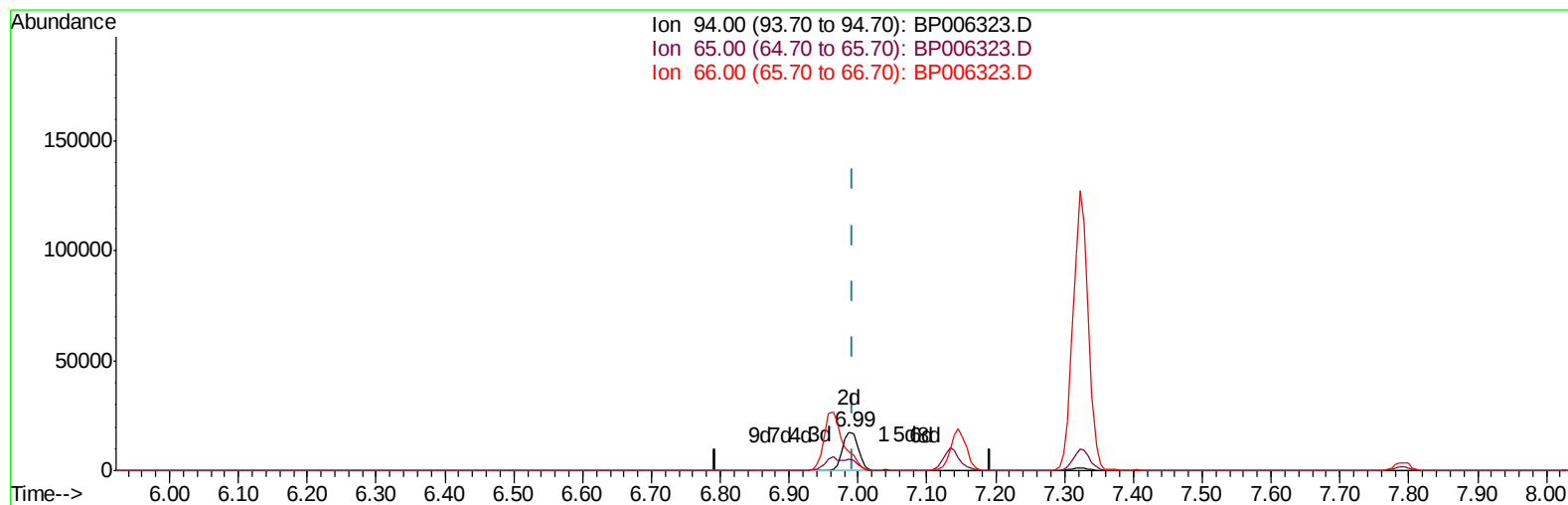
7.040min (+0.047) 0.01ng/ul

response 190

Ion	Exp%	Act%
94.00	100	100
65.00	28.20	25.77
66.00	36.90	31.90
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP072421\
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 Operator : CG/JU
 Sample : M2973-06
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Quant Time: Jul 26 01:10:08 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 24 04:29:12 2021
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TIC: BP006323.D

(8) Phenol

6.987min (-0.006) 1.00ng/ul m

response 28290

Ion	Exp%	Act%
94.00	100	100
65.00	28.20	28.47
66.00	36.90	49.15#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP072421\
 Data File : BP006323.D
 Acq On : 25 Jul 2021 01:21
 Operator : CG/JU
 Sample : M2973-06
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Quant Time: Jul 26 04:57:00 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 24 04:29:12 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	296123	20.00	ng/ul	0.00
20) Naphthalene-d8	10.57	136	1262419	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.42	164	720562	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.16	188	1394793	20.00	ng/ul	0.00
79) Chrysene-d12	21.26	240	1292331	20.00	ng/ul	0.00
88) Perylene-d12	23.61	264	1321284	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.31	96	48261	6.08	ng/uL	0.01
4) Pyridine-d5	3.71	84	686719	29.16	ng/ul	0.00
7) Phenol-d5	6.96	99	870378	31.30	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.13	67	556972	32.19	ng/ul	0.00
11) 2-Chlorophenol-d4	7.32	132	686877	33.24	ng/ul	0.00
15) 4-Methylphenol-d8	8.49	113	699815	32.00	ng/ul	0.00
21) Nitrobenzene-d5	8.95	128	321377	34.19	ng/ul	0.00
24) 2-Nitrophenol-d4	9.66	143	320317	35.69	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.19	165	566189	32.19	ng/ul	0.00
31) 4-Chloroaniline-d4	10.72	131	889624	30.72	ng/ul	0.00
46) Dimethylphthalate-d6	13.83	166	1812876	35.35	ng/ul	0.00
49) Acenaphthylene-d8	14.11	160	2193523	34.81	ng/ul	0.00
54) 4-Nitrophenol-d4	14.61	143	263573	25.88	ng/ul	0.00
60) Fluorene-d10	15.41	176	1495982	34.70	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.53	200	177770	22.86	ng/ul	0.00
73) Anthracene-d10	17.26	188	2235205	35.50	ng/ul	0.00
81) Pyrene-d10	19.50	212	2472104	37.19	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.46	264	2361615	35.17	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
72) Phenanthrene	17.21	178	94885	1.267	ng/ul	97

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

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