

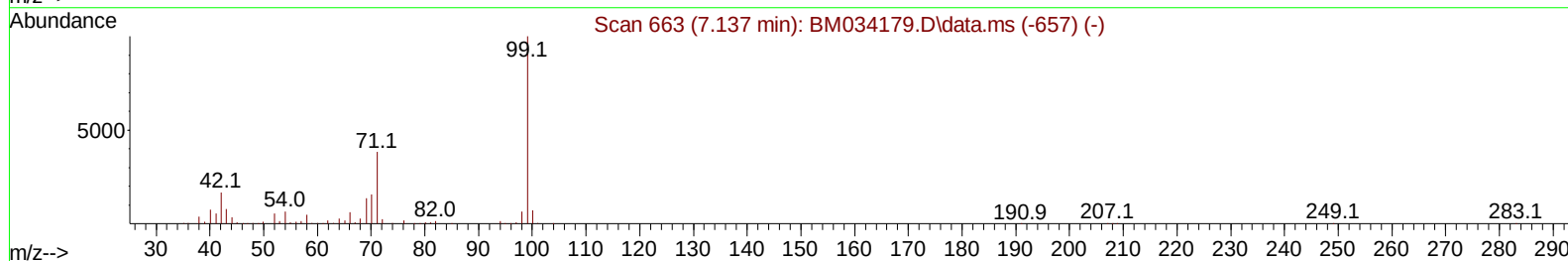
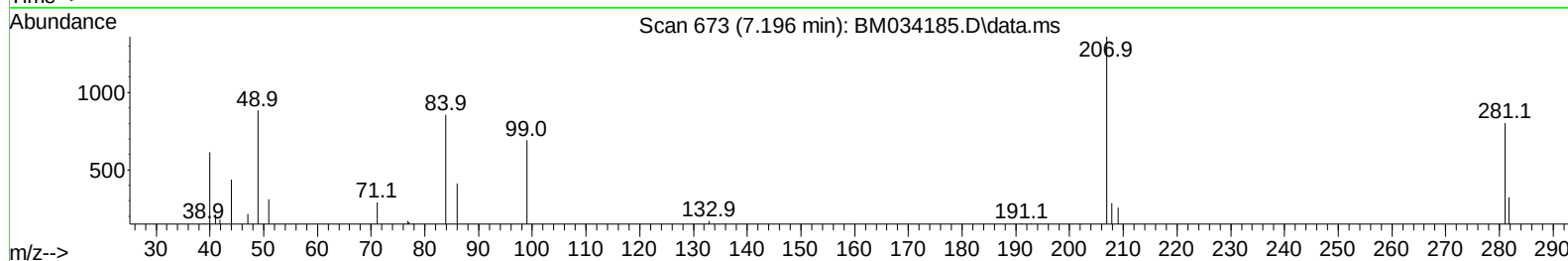
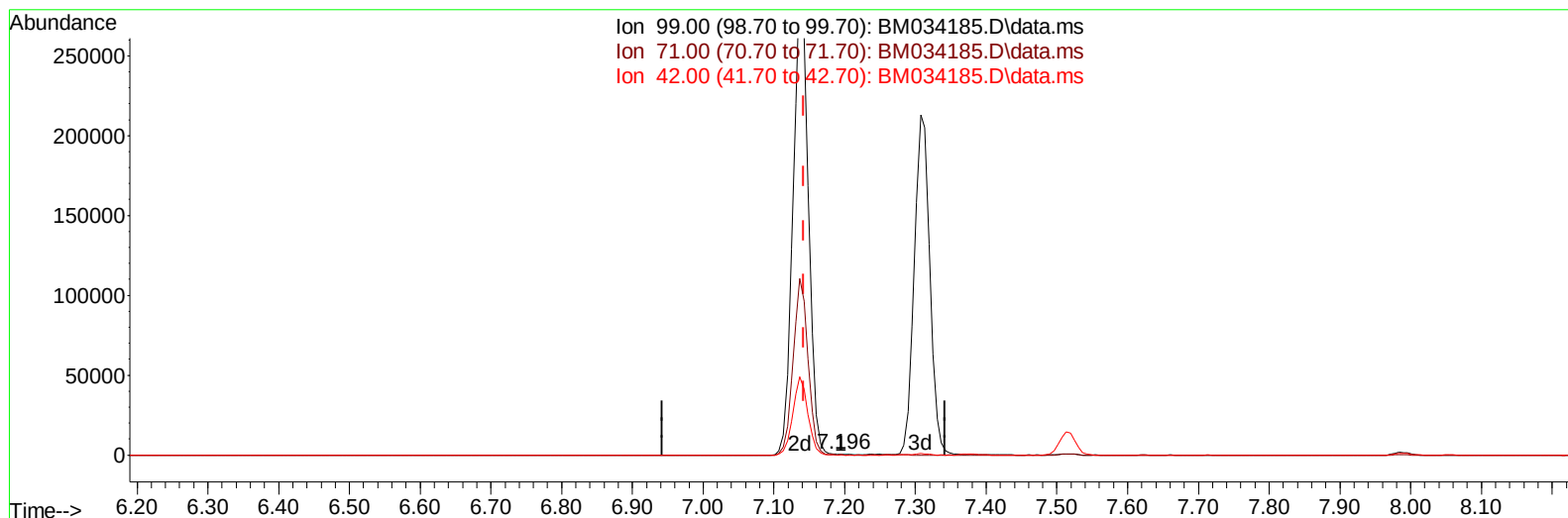
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
 Data File : BM034185.D
 Acq On : 09 Mar 2022 14:32
 Operator : CG/JU
 Sample : N1775-05
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBRV1

Manual IntegrationsAPPROVED

Quant Time: Mar 09 23:57:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 08 13:09:16 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/10/2022
 Supervised By :Yogesh Patel 03/11/2022



TIC: BM034185.D\data.ms

(7) Phenol-d5 (S)

7.196min (+ 0.053) 0.04 ng/ul

response 549

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	48.60	41.97
42.00	27.10	25.90
0.00	0.00	0.00

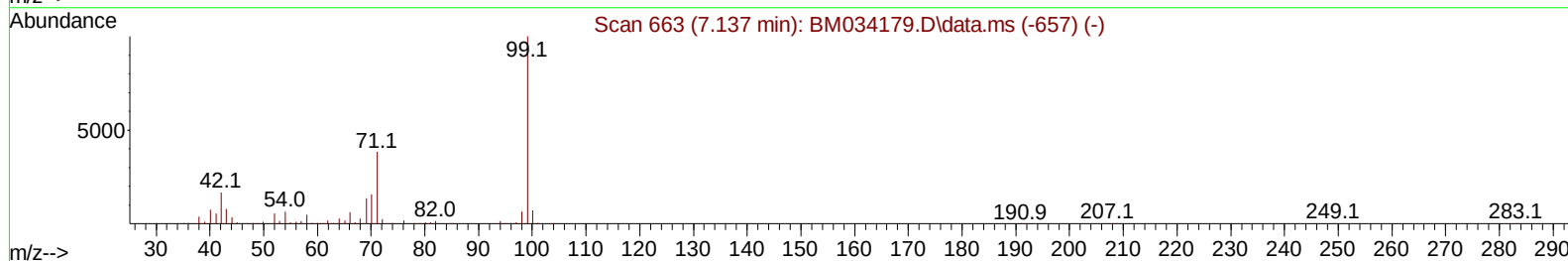
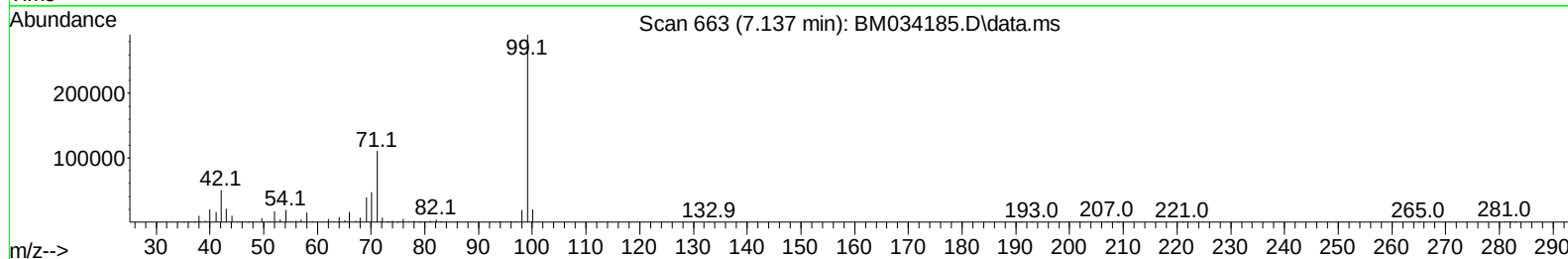
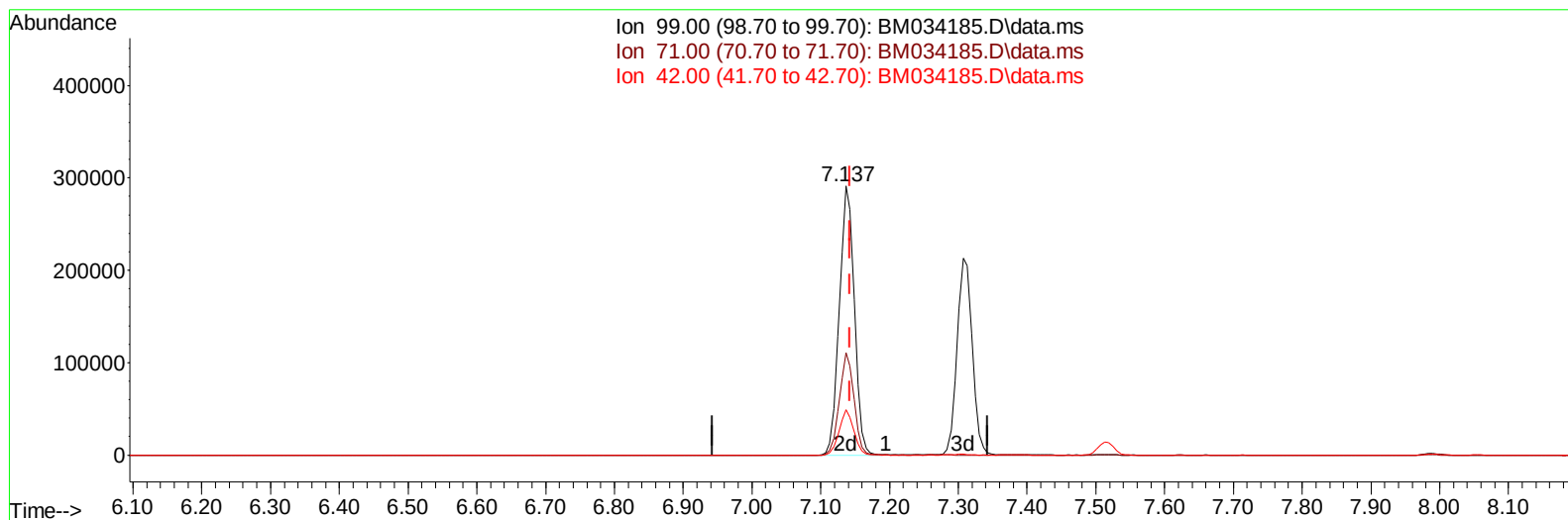
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(7) Phenol-d5 (S)

7.137min (-0.006) 33.06 ng/ul m

response 442328

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	48.60	37.99#
42.00	27.10	16.89#
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.990	152	170728	20.000	ng/ul	0.00
20) Naphthalene-d8	10.801	136	703328	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.624	164	487536	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.365	188	1029521	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.536	240	916175	20.000	ng/ul	0.00
88) Perylene-d12	23.983	264	858901	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.366	96	22949	6.182	ng/uL	0.00
4) Pyridine-d5	3.790	84	321221	29.727	ng/ul	0.00
7) Phenol-d5	7.137	99	442328m	33.065	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.307	67	247083	34.316	ng/ul	0.00
11) 2-Chlorophenol-d4	7.513	132	388201	35.007	ng/ul	0.00
15) 4-Methylphenol-d8	8.690	113	389383	34.738	ng/ul	0.00
21) Nitrobenzene-d5	9.148	128	192542	35.332	ng/ul	0.00
24) 2-Nitrophenol-d4	9.878	143	223415	37.747	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.419	165	421336	34.519	ng/ul	0.00
31) 4-Chloroaniline-d4	10.931	131	562310	33.718	ng/ul	0.00
46) Dimethylphthalate-d6	14.030	166	1432156	38.711	ng/ul	0.00
49) Acenaphthylene-d8	14.319	160	1687707	36.481	ng/ul	0.00
54) 4-Nitrophenol-d4	14.801	143	200053	29.061	ng/ul	0.00
60) Fluorene-d10	15.613	176	1217469	36.976	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.718	200	191534	28.829	ng/ul	0.00
73) Anthracene-d10	17.465	188	1939046	38.235	ng/ul	0.00
81) Pyrene-d10	19.753	212	2180301	43.295	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.824	264	1792575	39.343	ng/ul	0.00

Target Compounds Qvalue

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

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