

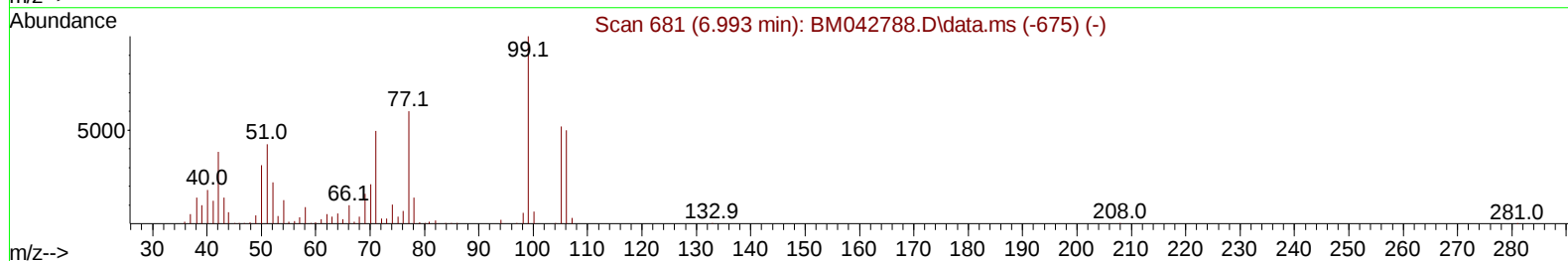
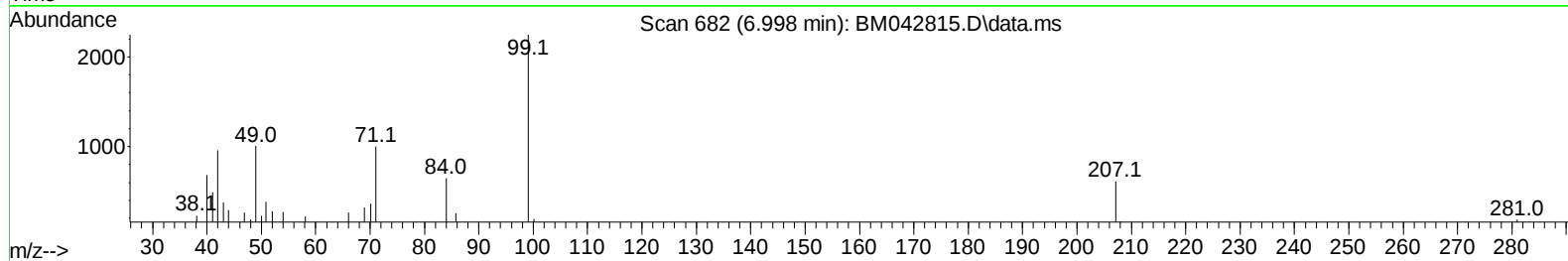
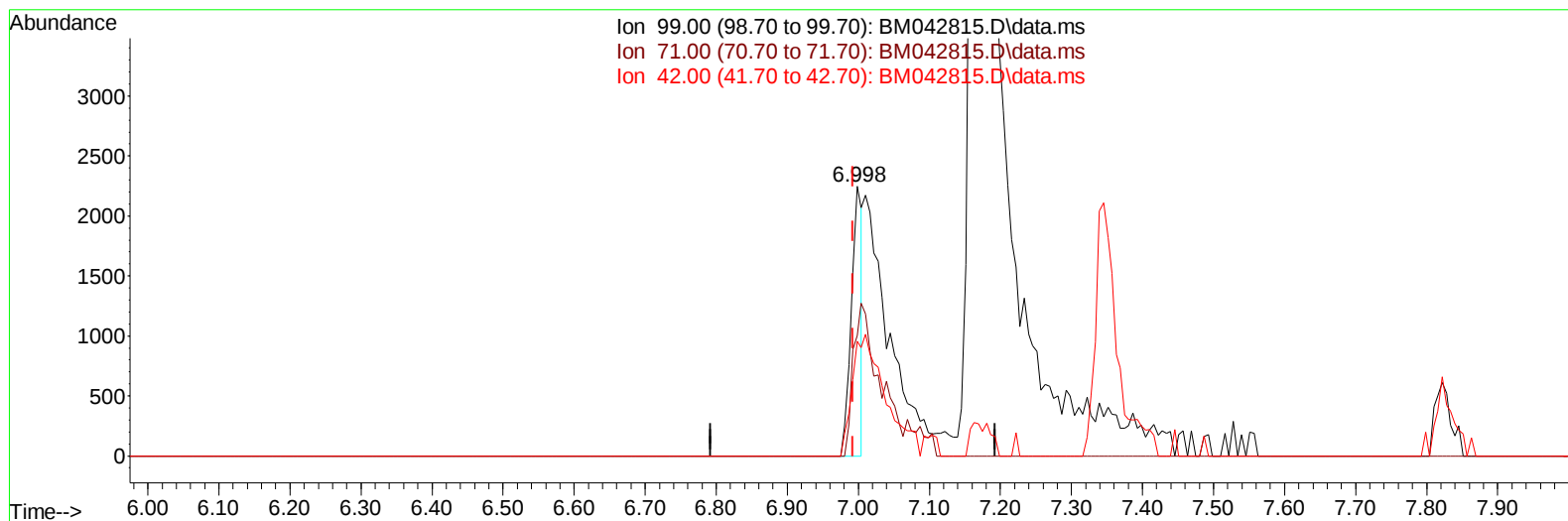
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
Data File : BM042815.D
Acq On : 16 Nov 2023 20:56
Operator : MA/JU
Sample : 05268-06
Misc :
ALS Vial : 57 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCKC8

Manual IntegrationsAPPROVED

Quant Time: Nov 16 23:56:09 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 15 22:28:07 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/17/2023
Supervised By :mohammad ahmed 11/19/2023



TIC: BM042815.D\data.ms

(7) Phenol-d5 (S)

6.998min (+ 0.006) 1.05 ng/u1

response 2420

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	43.50	44.57
42.00	32.20	42.65#
0.00	0.00	0.00

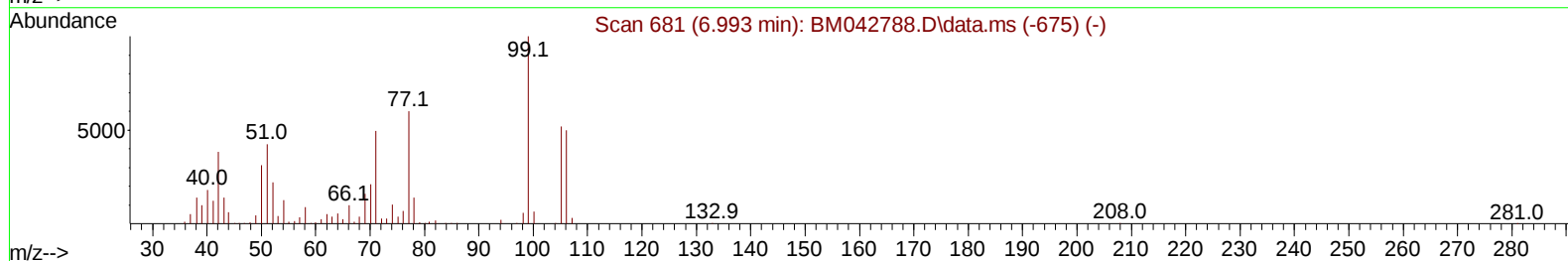
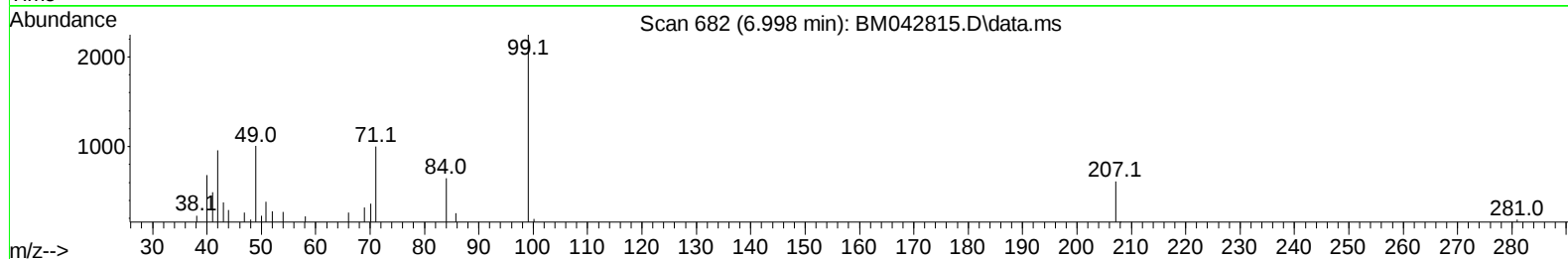
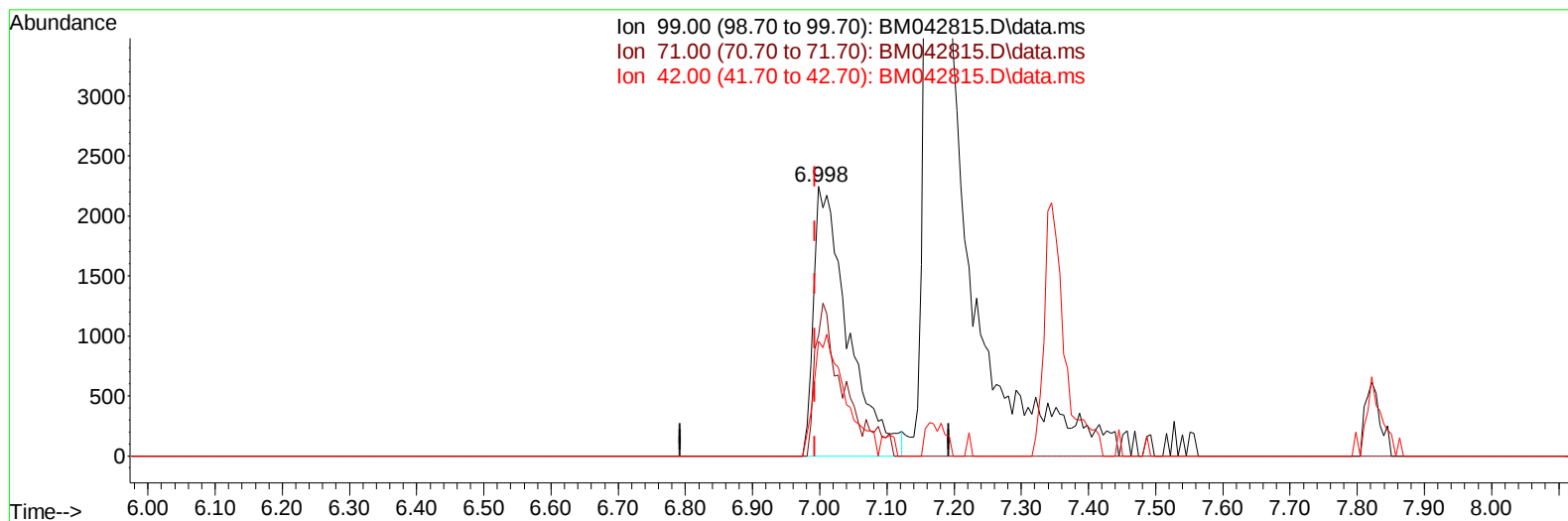
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TIC: BM042815.D\data.ms

(7) Phenol-d5 (S)

6.998min (+ 0.006) 3.46 ng/ul m

response 7960

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	43.50	44.57
42.00	32.20	42.65#
0.00	0.00	0.00

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Quant Time: Nov 17 00:34:54 2023
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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 22:28:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	23639	20.000	ng/ul	0.00
20) Naphthalene-d8	10.633	136	90607	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.480	164	56961	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.233	188	123222	20.000	ng/ul	0.00
79) Chrysene-d12	21.415	240	128496	20.000	ng/ul	# 0.00
88) Perylene-d12	23.780	264	170169	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	1583	2.101	ng/uL	0.00
4) Pyridine-d5	3.622	84	21754	11.609	ng/ul	0.00
7) Phenol-d5	6.998	99	7960m	3.460	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.169	67	35377	21.883	ng/ul	0.00
11) 2-Chlorophenol-d4	7.345	132	24713	13.921	ng/ul	0.00
15) 4-Methylphenol-d8	8.557	113	13412	7.404	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	14974	18.850	ng/ul	0.00
24) 2-Nitrophenol-d4	9.739	143	13503	14.530	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.269	165	25445	14.726	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	34659	15.798	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166	125684	23.314	ng/ul	0.00
49) Acenaphthylene-d8	14.180	160	126226	21.883	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0	0.000	ng/ul	
60) Fluorene-d10	15.474	176	102998	23.072	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.627	200	9429	10.102	ng/ul	0.00
73) Anthracene-d10	17.333	188	167600	24.270	ng/ul	0.00
81) Pyrene-d10	19.627	212	233360	27.733	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.621	264	260308	25.574	ng/ul	0.00

Target Compounds Qvalue

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

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