

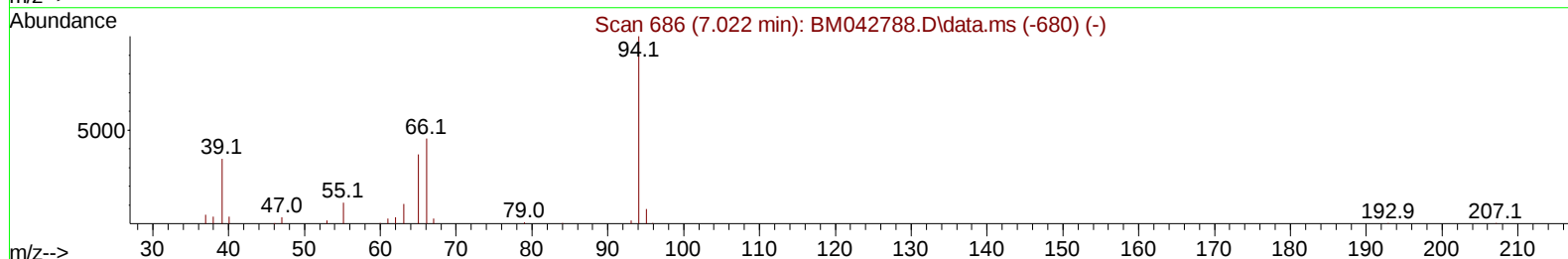
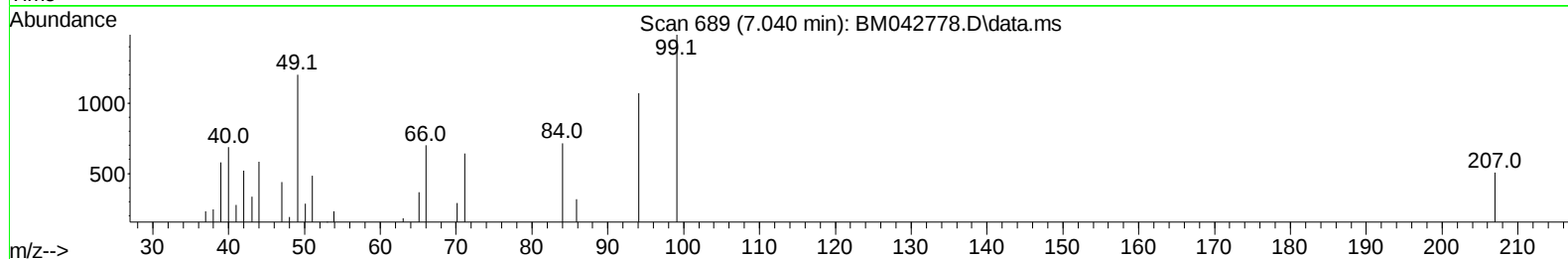
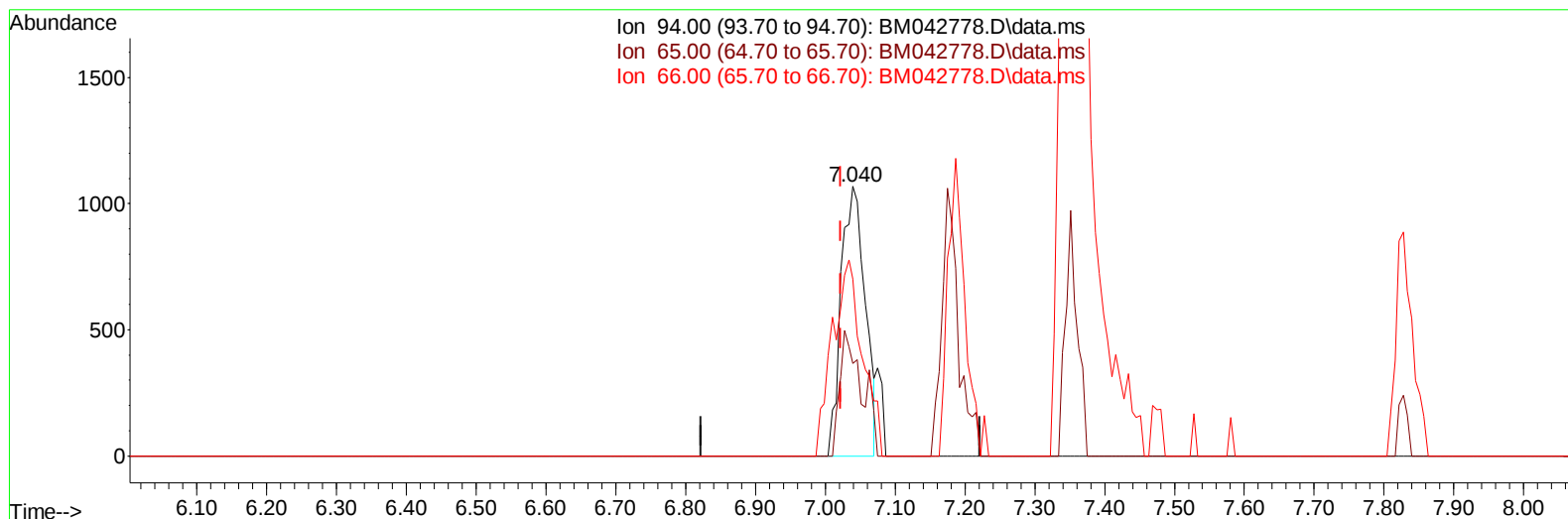
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
Data File : BM042778.D
Acq On : 15 Nov 2023 20:54
Operator : MA/JU
Sample : 05057-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A49F7

Manual IntegrationsAPPROVED

Quant Time: Nov 17 04:02:25 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 11/16/2023
Supervised By :mohammad ahmed 11/19/2023



TIC: BM042778.D\data.ms

(8) Phenol

7.040min (+ 0.017) 1.27 ng/u1

response 2535

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	37.90	34.36
66.00	49.90	65.54#
0.00	0.00	0.00

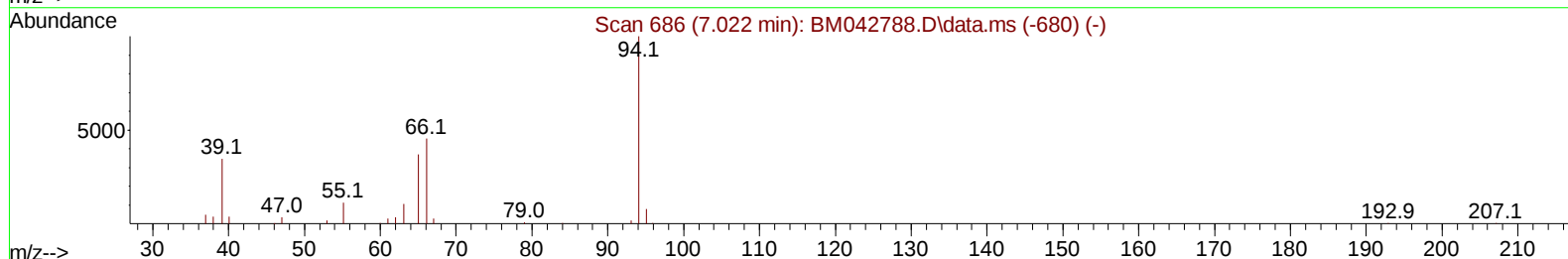
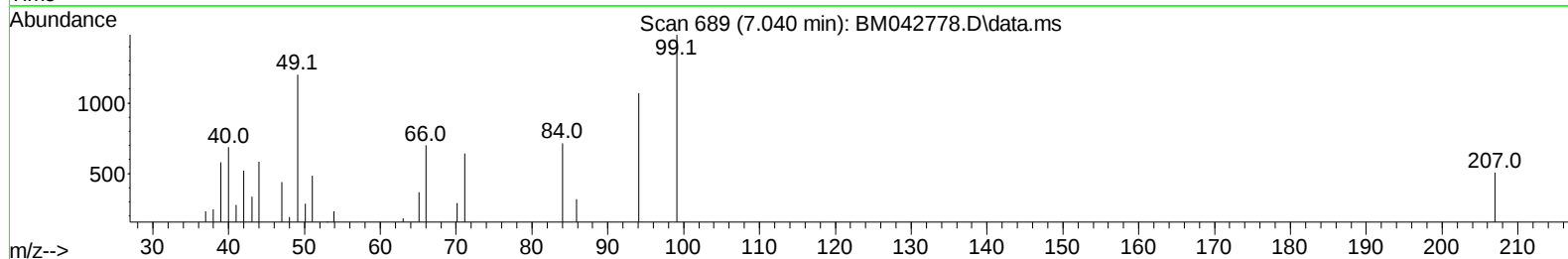
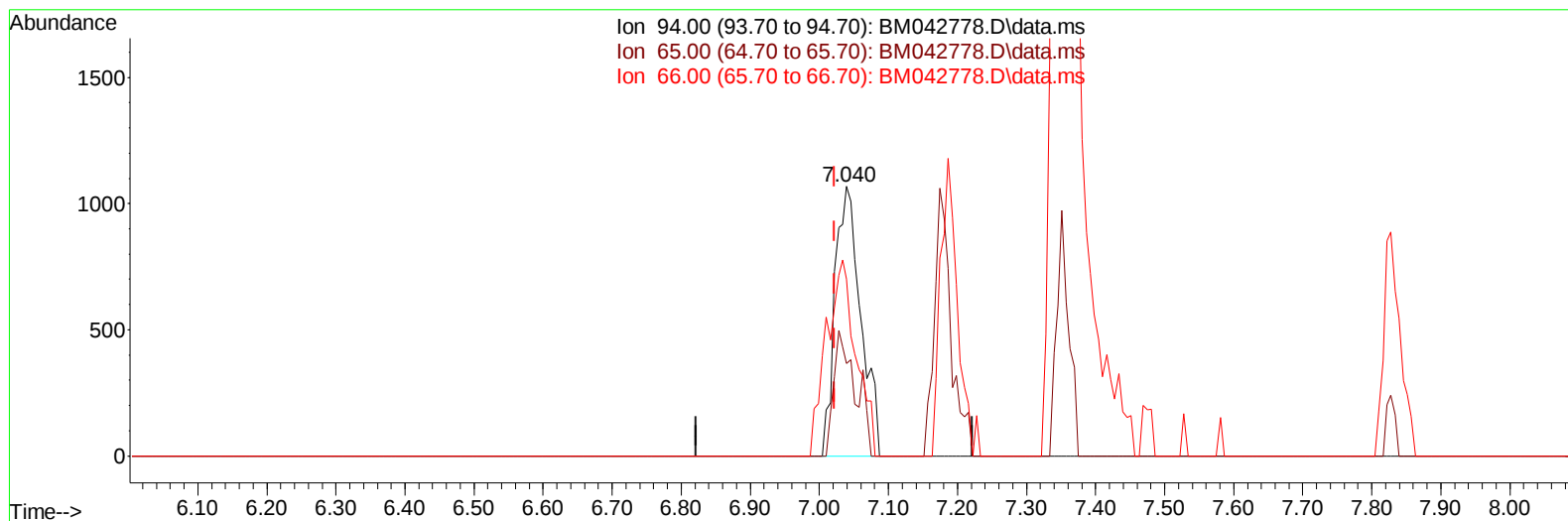
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Manual IntegrationsAPPROVED

Quant Time: Nov 15 22:49:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
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TIC: BM042778.D\data.ms

(8) Phenol

7.040min (+ 0.017) 1.38 ng/ul m

response 2758

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	37.90	34.36
66.00	49.90	65.54#
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	7.828	152	20693	20.000	ng/ul	0.00
20) Naphthalene-d8	10.639	136	87158	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.480	164	56018	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.233	188	125377	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.421	240	126531	20.000	ng/ul	# 0.00
88) Perylene-d12	23.785	264	156209	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	2627	3.983	ng/uL	0.00
4) Pyridine-d5	3.634	84	8854	5.397	ng/ul	0.02
7) Phenol-d5	7.004	99	9825	4.878	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.175	67	43702	30.881	ng/ul	0.00
11) 2-Chlorophenol-d4	7.351	132	32275	20.769	ng/ul	0.00
15) 4-Methylphenol-d8	8.557	113	18362	11.579	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	19782	25.887	ng/ul	0.00
24) 2-Nitrophenol-d4	9.745	143	19844	22.199	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.269	165	35158	21.152	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	40423	19.154	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166	165496	31.216	ng/ul	0.00
49) Acenaphthylene-d8	14.186	160	163505	28.824	ng/ul	0.00
54) 4-Nitrophenol-d4	14.804	143	227	0.379	ng/ul	0.06
60) Fluorene-d10	15.474	176	135742	30.919	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.627	200	16671	17.554	ng/ul	0.00
73) Anthracene-d10	17.339	188	210335	29.935	ng/ul	0.00
81) Pyrene-d10	19.633	212	310335	37.454	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.627	264	310801	33.263	ng/ul	0.00
Target Compounds						
8) Phenol	7.040	94	2758m	1.377	ng/ul	Qvalue

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

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