Quantitation Report (Qedit)

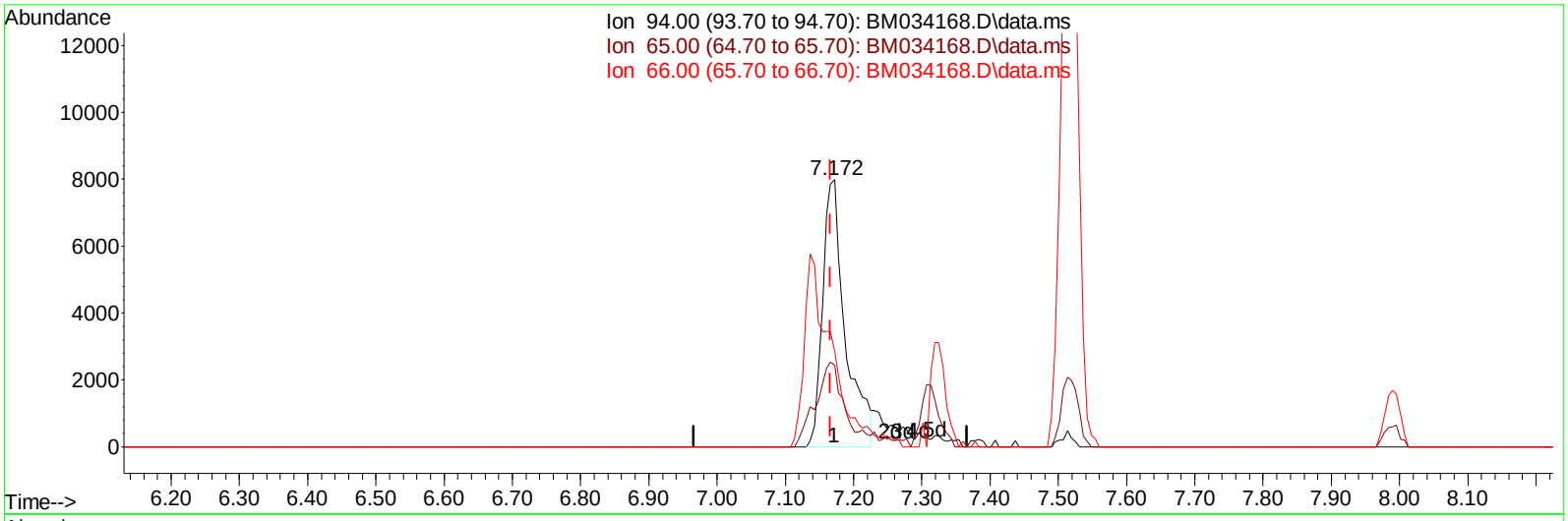
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
Data File : BM034168.D
Acq On : 09 Mar 2022 03:07
Operator : CG/JU
Sample : N1733-23
Misc :
ALS Vial : 27 Sample Multiplier: 1

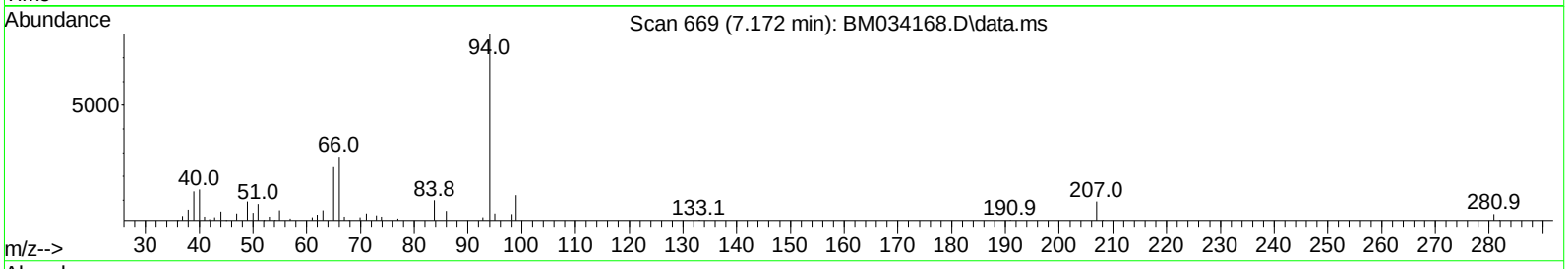
Quant Time: Mar 09 03:47:17 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Mar 08 13:09:16 2022
Response via : Initial Calibration

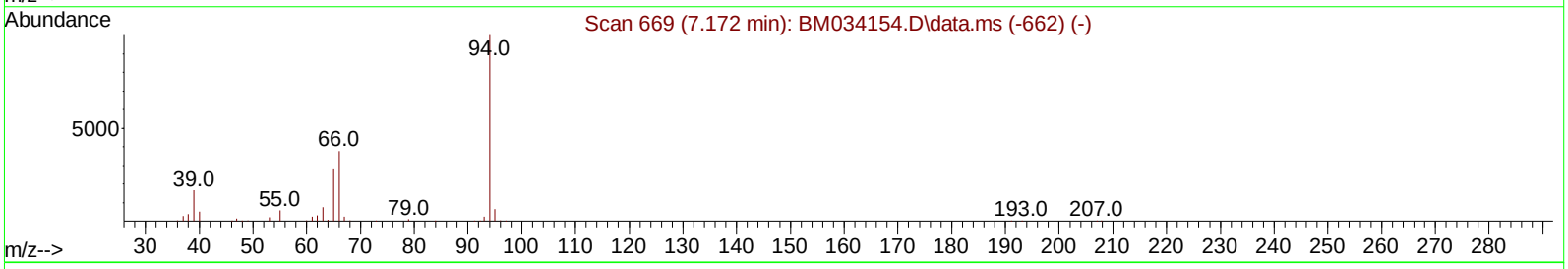
Instrument :
BNA_M
ClientSampleId :
DBRU1

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 03/09/2022
Supervised By :Yogesh Patel 03/09/2022







TIC: BM034168.D\data.ms

(8) Phenol

7.172min (+ 0.006) 1.68 ng/ul

response 18469

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	39.60	30.65#
66.00	56.70	35.64#
0.00	0.00	0.00

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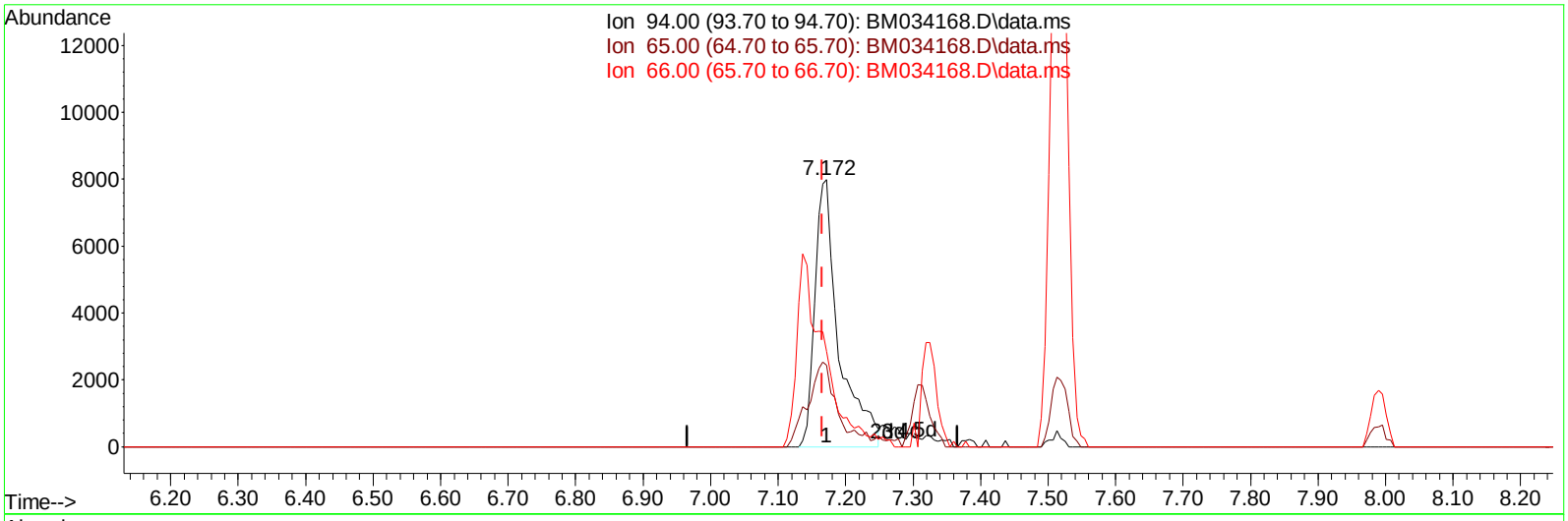

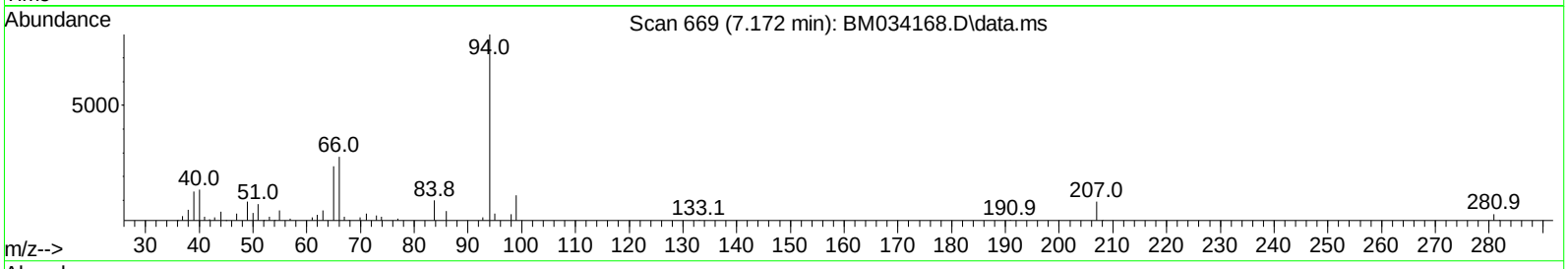

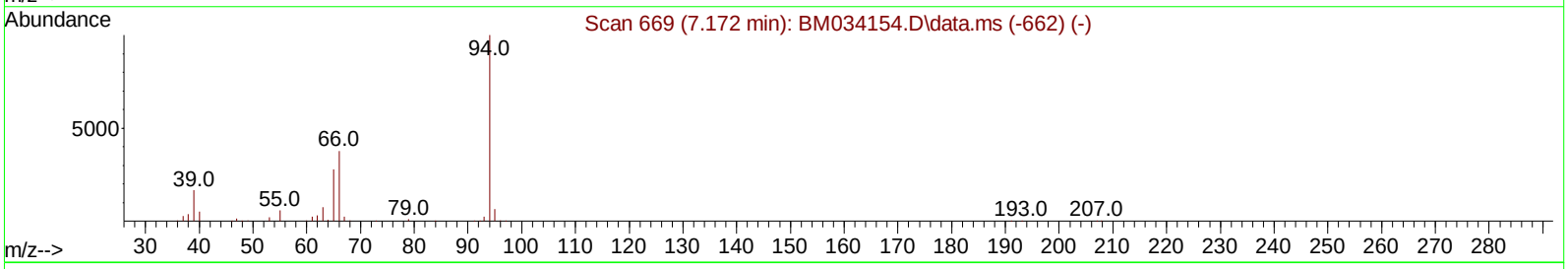

TIC: BM034168.D\data.ms

(8) Phenol

7.172min (+ 0.006) 1.79 ng/ul m

response 19663

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	39.60	30.65#
66.00	56.70	35.64#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.990	152	138539	20.000	ng/ul	0.00
20) Naphthalene-d8	10.807	136	498021	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.624	164	311017	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.365	188	628859	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.536	240	667724	20.000	ng/ul	# 0.00
88) Perylene-d12	23.983	264	776147	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	9918	3.293	ng/uL	0.00
4) Pyridine-d5	3.796	84	53220	6.070	ng/ul	0.00
7) Phenol-d5	7.137	99	158353	14.588	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.307	67	91429	15.648	ng/ul	0.00
11) 2-Chlorophenol-d4	7.519	132	144473	16.055	ng/ul	0.00
15) 4-Methylphenol-d8	8.690	113	119234	13.109	ng/ul	0.00
21) Nitrobenzene-d5	9.148	128	65956	17.093	ng/ul	0.00
24) 2-Nitrophenol-d4	9.878	143	71136	16.973	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.419	165	134305	15.539	ng/ul	0.00
31) 4-Chloroaniline-d4	10.931	131	154795	13.109	ng/ul	0.00
46) Dimethylphthalate-d6	14.030	166	411640	17.442	ng/ul	0.00
49) Acenaphthylene-d8	14.319	160	482953	16.364	ng/ul	0.00
54) 4-Nitrophenol-d4	14.801	143	50539	11.508	ng/ul	0.00
60) Fluorene-d10	15.613	176	348144	16.574	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.718	200	49345	12.159	ng/ul	0.00
73) Anthracene-d10	17.465	188	539963	17.431	ng/ul	0.00
81) Pyrene-d10	19.748	212	653284	17.799	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.818	264	728924	17.704	ng/ul	-0.01
Target Compounds						Qvalue
8) Phenol	7.172	94	19663m	1.791	ng/ul	

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

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