Quantitation Report (Qedit)

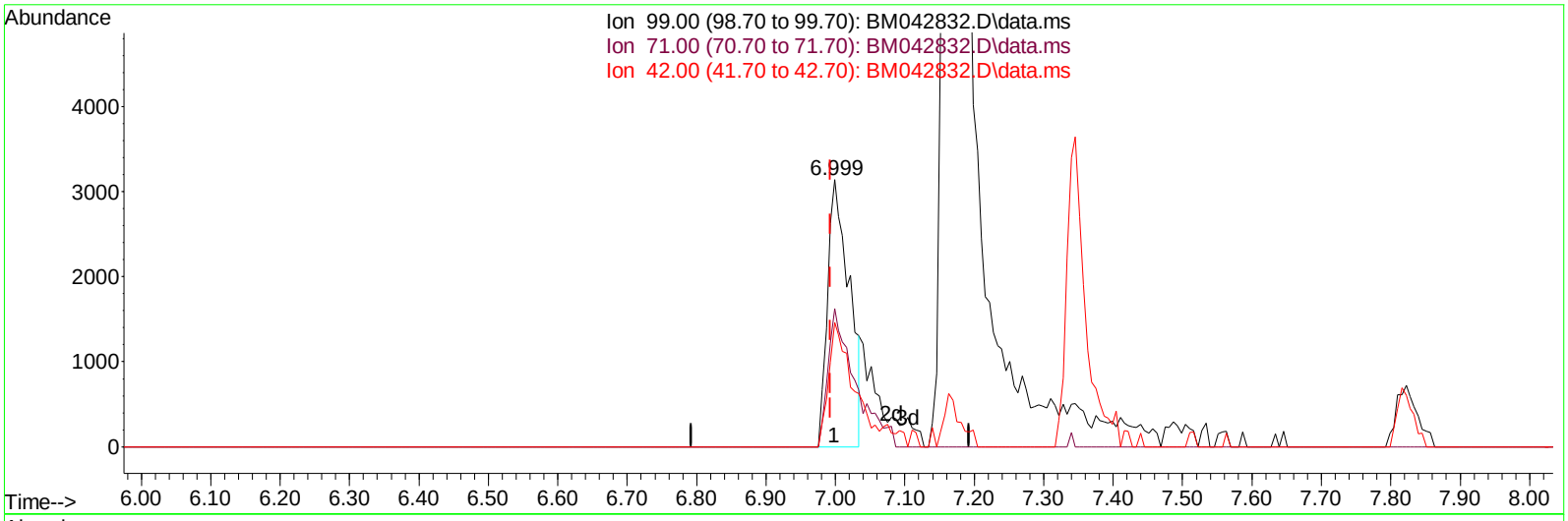
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
Data File : BM042832.D
Acq On : 17 Nov 2023 07:46
Operator : MA/JU
Sample : 05268-18
Misc :
ALS Vial : 75 Sample Multiplier: 1

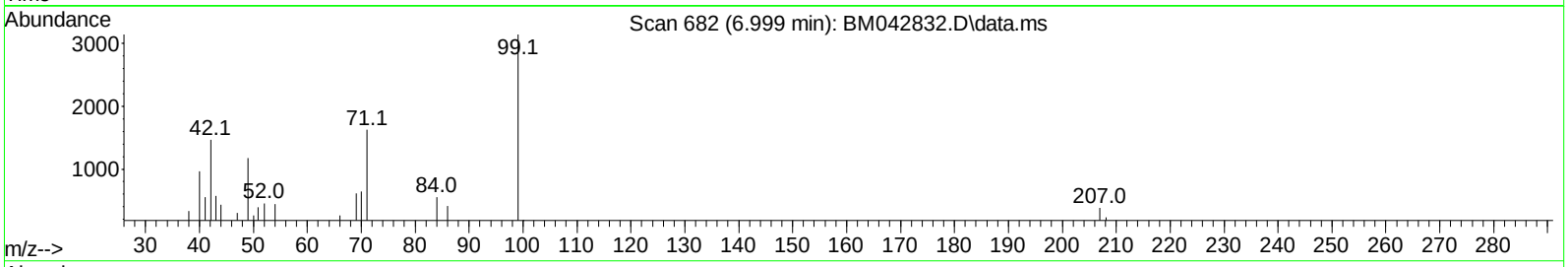
Quant Time: Nov 17 11:39:11 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 15 22:28:07 2023
Response via : Initial Calibration

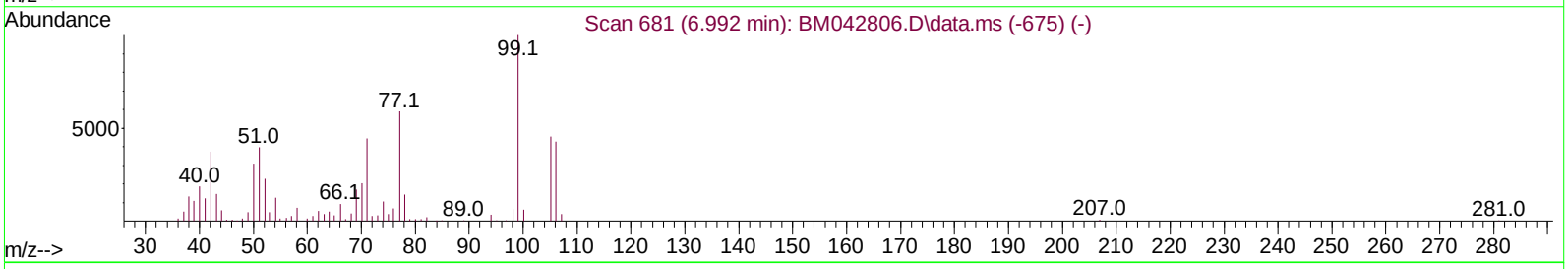
Instrument :
BNA_M
ClientSampleId :
DCKE6

Manual IntegrationsAPPROVED

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







TIC: BM042832.D\data.ms

(7) Phenol-d5 (S)

6.999min (+ 0.006) 2.75 ng/ul

response 6909

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	43.50	51.72
42.00	32.20	46.69#
0.00	0.00	0.00

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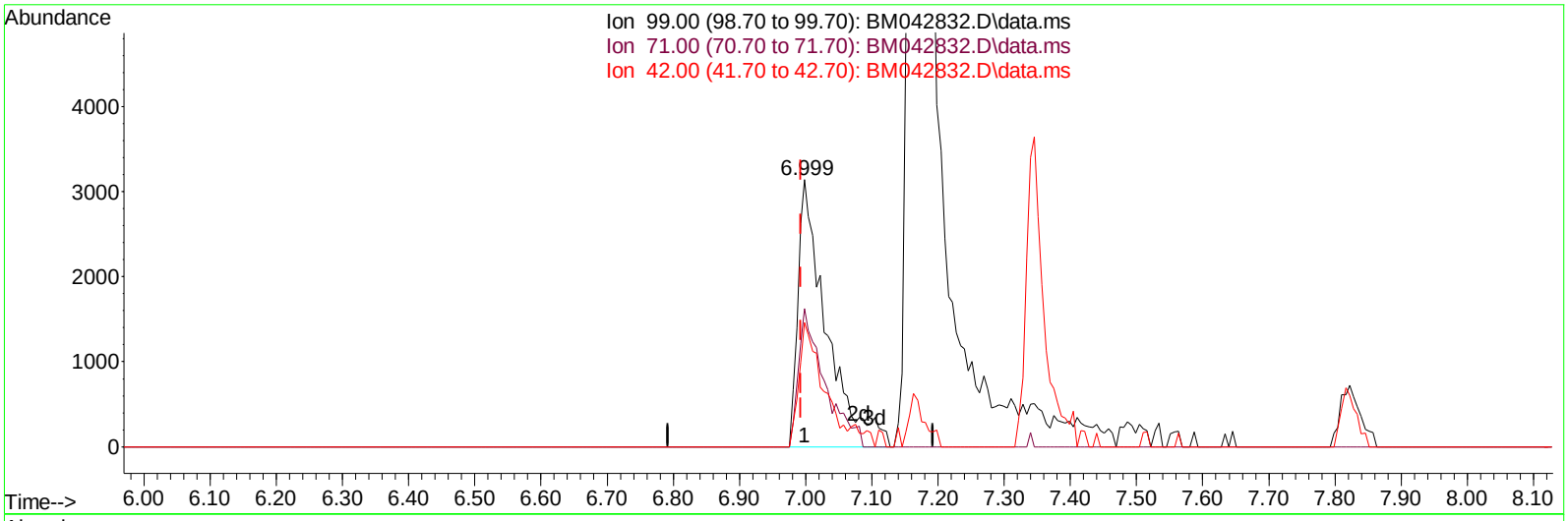

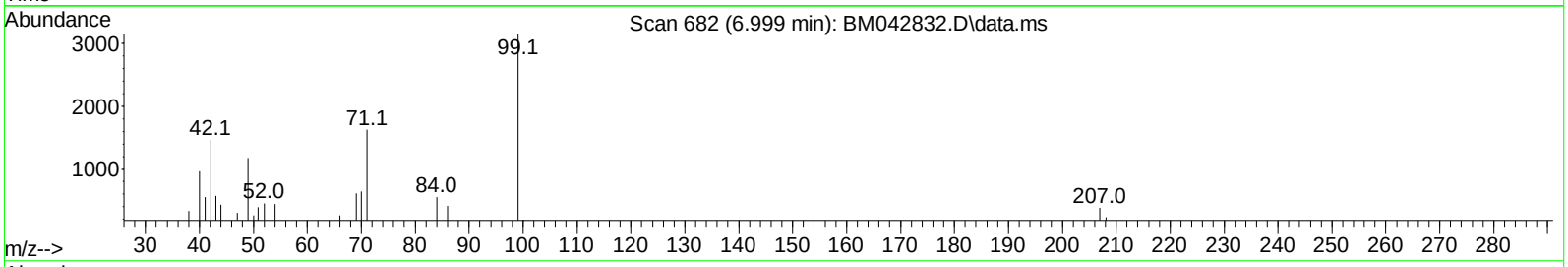

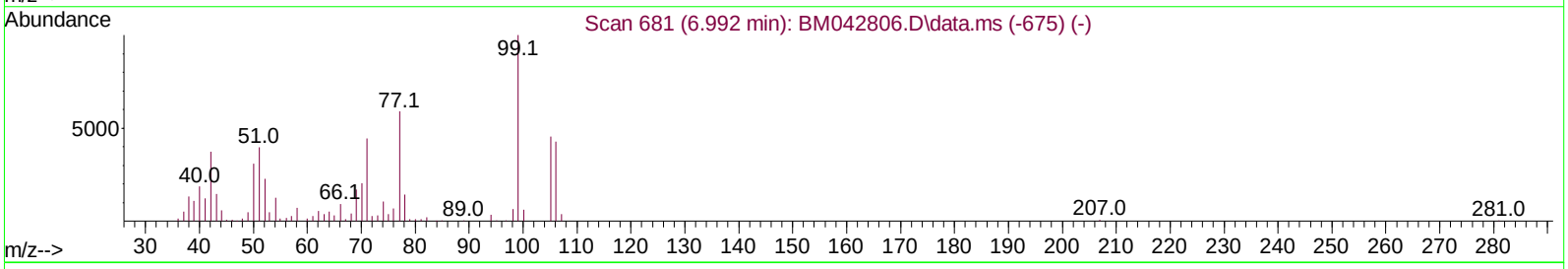

TIC: BM042832.D\data.ms

(7) Phenol-d5 (S)

6.999min (+ 0.006) 3.72 ng/ul m

response 9357

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	43.50	51.72
42.00	32.20	46.69#
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	7.822	152	25860	20.000	ng/ul	0.00
20) Naphthalene-d8	10.634	136	105815	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.475	164	63260	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.233	188	140301	20.000	ng/ul	0.00
79) Chrysene-d12	21.415	240	136018	20.000	ng/ul	0.00
88) Perylene-d12	23.774	264	170362	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	2523	3.061	ng/uL	0.00
4) Pyridine-d5	3.628	84	13089	6.385	ng/ul	0.01
7) Phenol-d5	6.999	99	9357m	3.718	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.169	67	57621	32.581	ng/ul	0.00
11) 2-Chlorophenol-d4	7.346	132	37862	19.496	ng/ul	0.00
15) 4-Methylphenol-d8	8.551	113	17941	9.053	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	25266	27.234	ng/ul	0.00
24) 2-Nitrophenol-d4	9.734	143	25496	23.492	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.263	165	42306	20.965	ng/ul	0.00
31) 4-Chloroaniline-d4	10.816	131	37585	14.669	ng/ul	0.00
46) Dimethylphthalate-d6	13.898	166	180887	30.213	ng/ul	0.00
49) Acenaphthylene-d8	14.175	160	186378	29.094	ng/ul	-0.01
54) 4-Nitrophenol-d4	0.000	143	0	0.000	ng/ul	
60) Fluorene-d10	15.474	176	159486	32.168	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.621	200	18108	17.039	ng/ul	0.00
73) Anthracene-d10	17.327	188	257358	32.731	ng/ul	-0.01
81) Pyrene-d10	19.627	212	338723	38.029	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.621	264	345563	33.911	ng/ul	0.00

Target Compounds Qvalue

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

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