Quantitation Report (Qedit)

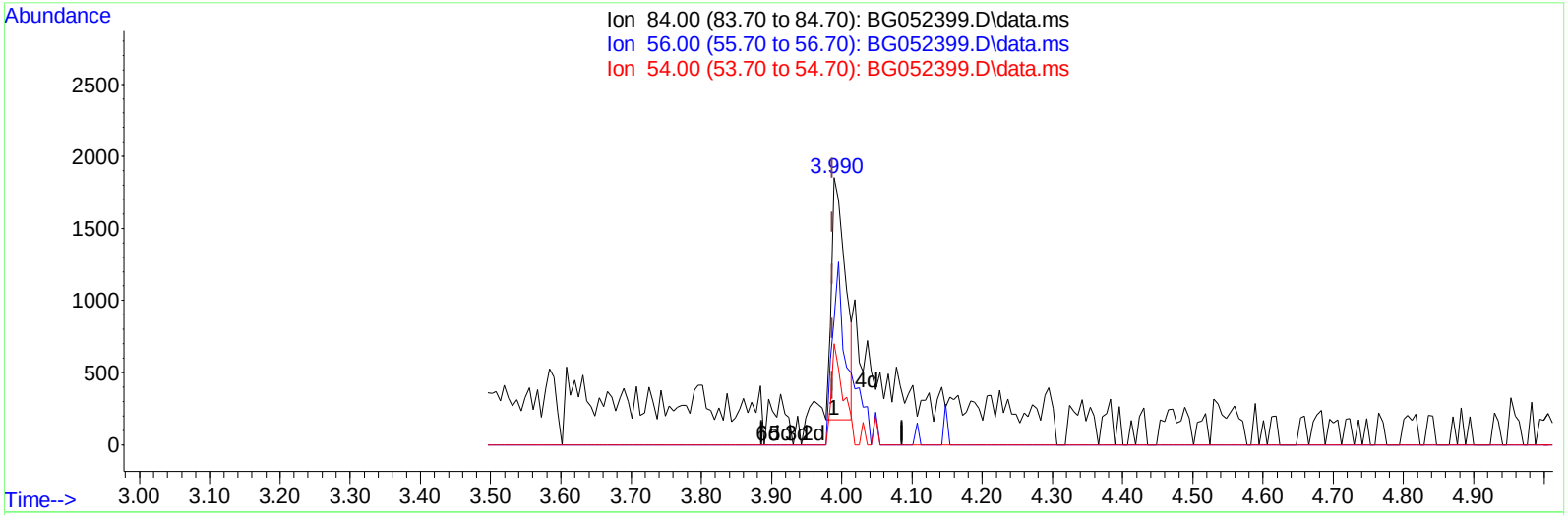
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021522\
Data File : BG052399.D
Acq On : 15 Feb 2022 19:30
Operator : CG/JU
Sample : N1478-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

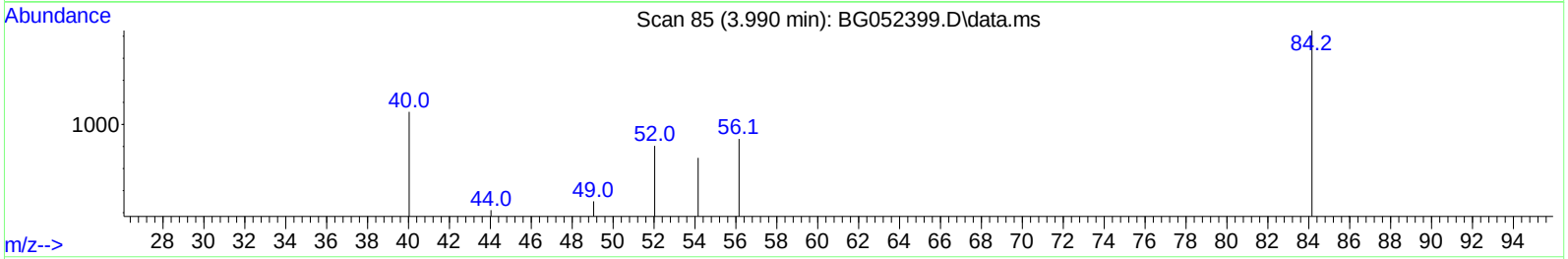
Quant Time: Feb 15 21:26:18 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG013122.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jan 31 18:07:10 2022
Response via : Initial Calibration

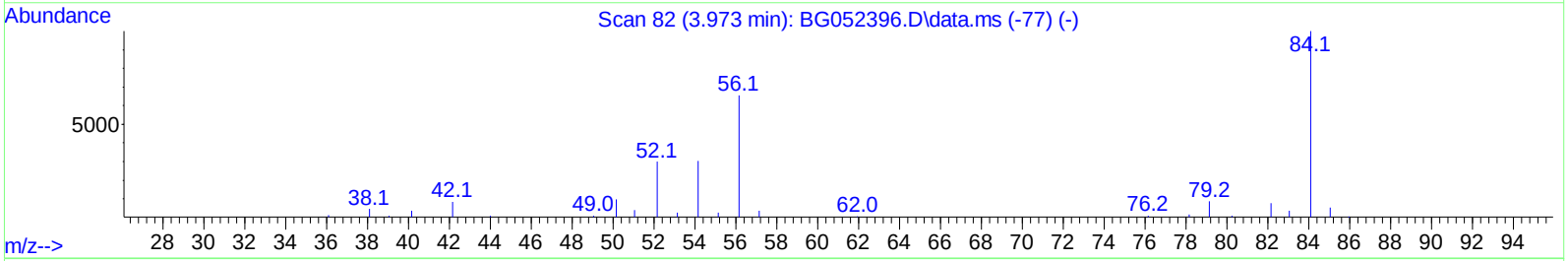
Instrument :
BNA_G
ClientSampleId :
GBCT9

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 02/16/2022
Supervised By :Jagrut Upadhyay 02/16/2022







TIC: BG052399.D\data.ms

(4) Pyridine-d5 (S)

3.990min (+ 0.003) 1.34 ng/ul

response 2319

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	47.00#
54.00	31.50	37.82#
0.00	0.00	0.00

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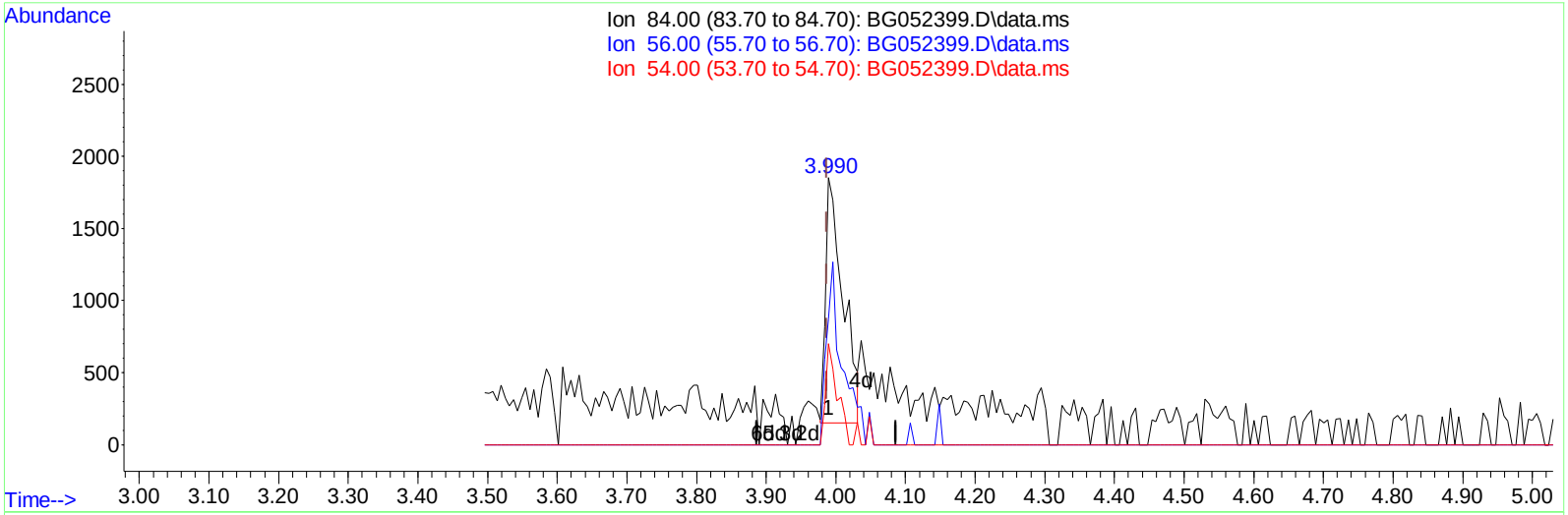

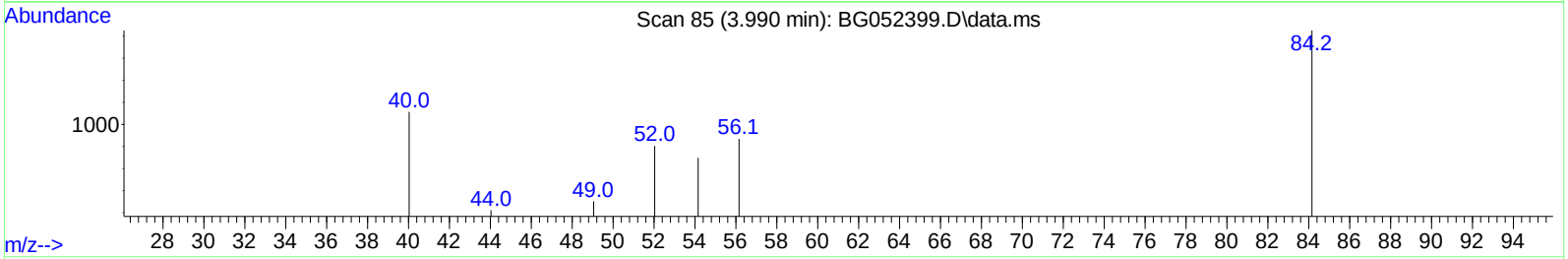

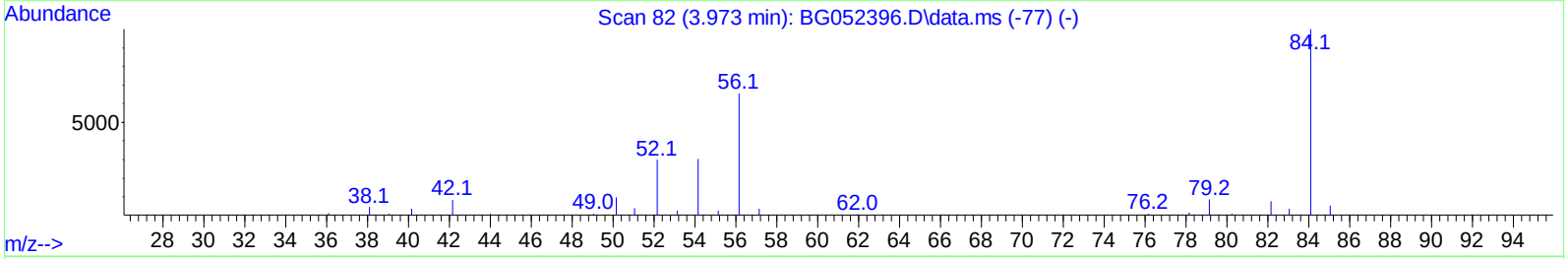

TIC: BG052399.D\data.ms

(4) Pyridine-d5 (S)

3.990min (+ 0.003) 1.69 ng/ul m

response 2936

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	47.00#
54.00	31.50	37.82#
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	8.219	152	22497	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.056	136	97209	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.863	164	71706	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.612	188	180884	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.924	240	202159	20.000	ng/ul	#-0.02
88) Perylene-d12	25.390	264	200759	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.543	96	2837	5.087	ng/uL	-0.02
4) Pyridine-d5	3.990	84	2936m	1.693	ng/ul	0.00
7) Phenol-d5	7.367	99	13906	6.796	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.526	67	39418	31.707	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.749	132	35550	24.931	ng/ul	-0.01
15) 4-Methylphenol-d8	8.924	113	25427	16.374	ng/ul	-0.01
21) Nitrobenzene-d5	9.394	128	25219	31.862	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.122	143	26533	30.677	ng/ul	-0.02
28) 2,4-Dichlorophenol-d3	10.675	165	45324	26.881	ng/ul	-0.01
31) 4-Chloroaniline-d4	0.000	131	0	0.000	ng/ul	
46) Dimethylphthalate-d6	14.252	166	199792	35.447	ng/ul	-0.01
49) Acenaphthylene-d8	14.563	160	218755	31.138	ng/ul	-0.01
54) 4-Nitrophenol-d4	15.063	143	6533	7.653	ng/ul	0.00
60) Fluorene-d10	15.856	176	180309	36.515	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.967	200	34322	30.735	ng/ul	-0.01
73) Anthracene-d10	17.712	188	334560	39.233	ng/ul	-0.01
81) Pyrene-d10	19.991	212	433663	35.793	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.155	264	403136	37.798	ng/ul	-0.03

Target Compounds Qvalue

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

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