Quantitation Report (Qedit)

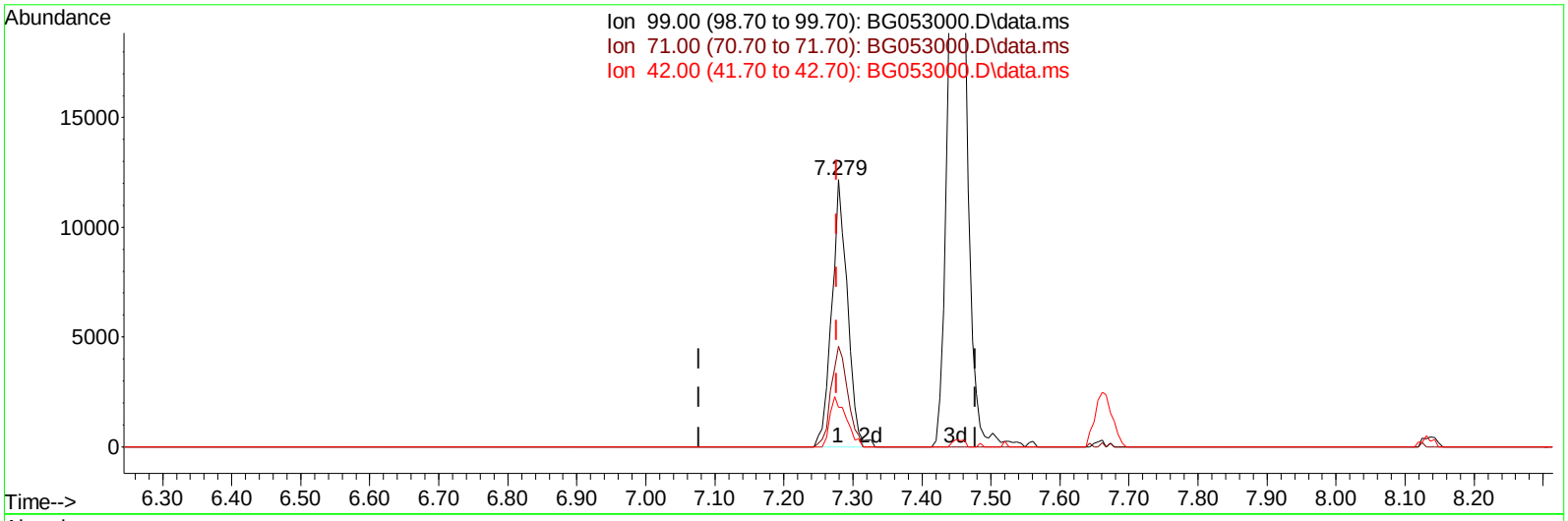
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG033022\
Data File : BG053000.D
Acq On : 1 Apr 2022 16:55
Operator : CG/JU
Sample : N2166-01
Misc :
ALS Vial : 60 Sample Multiplier: 1

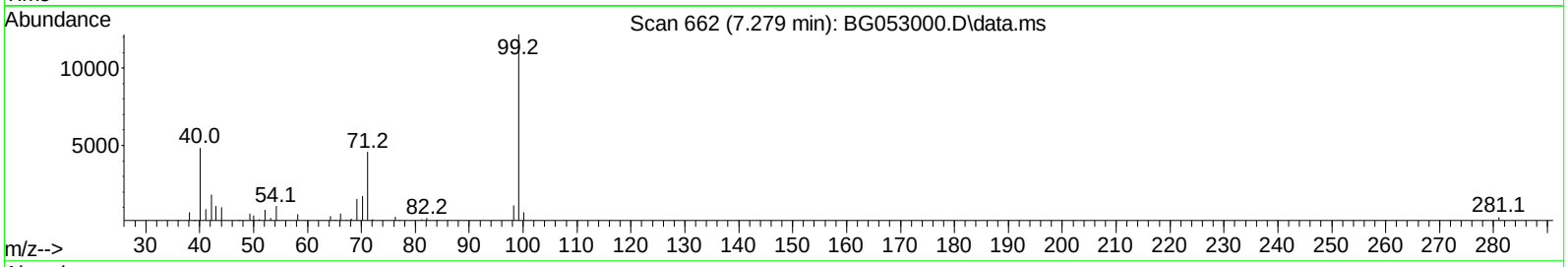
Quant Time: Apr 02 00:37:54 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 31 00:50:49 2022
Response via : Initial Calibration

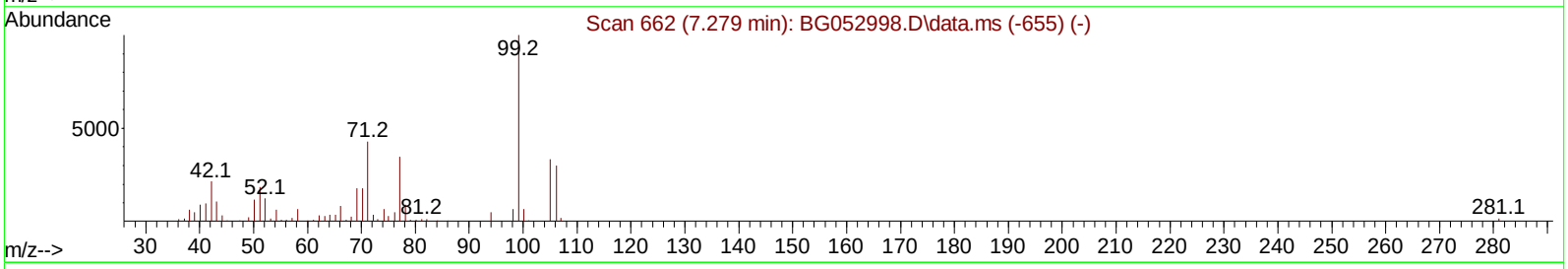
Instrument :
BNA_G
ClientSampleId :
C0R39

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/04/2022
Supervised By :Yogesh Patel 04/06/2022







TIC: BG053000.D\data.ms

(7) Phenol-d5 (S)

7.279min (+ 0.002) 7.41 ng/ul

response 19156

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	36.40	37.64
42.00	19.00	14.87#
0.00	0.00	0.00

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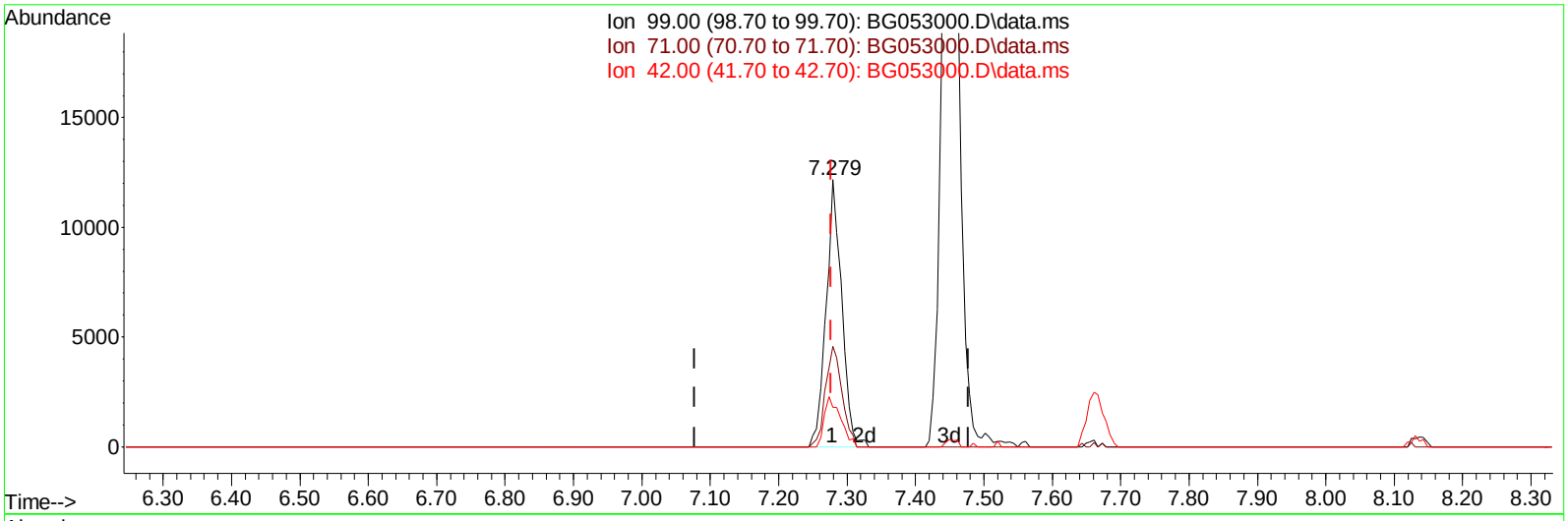

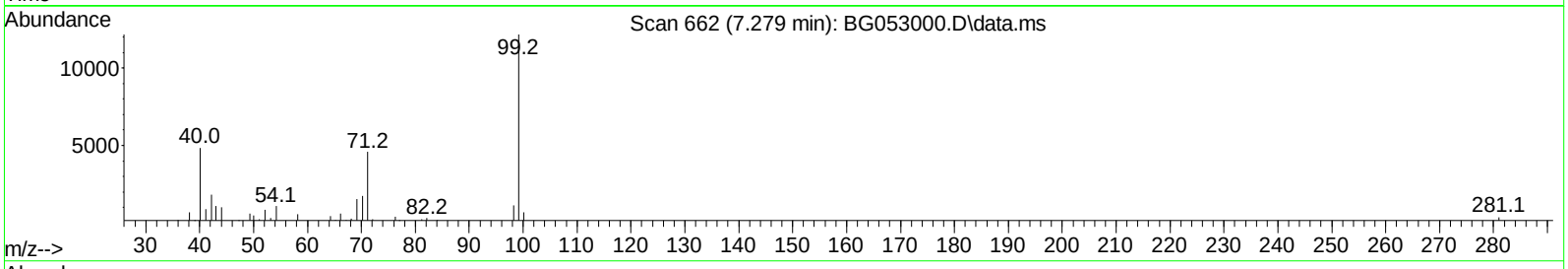

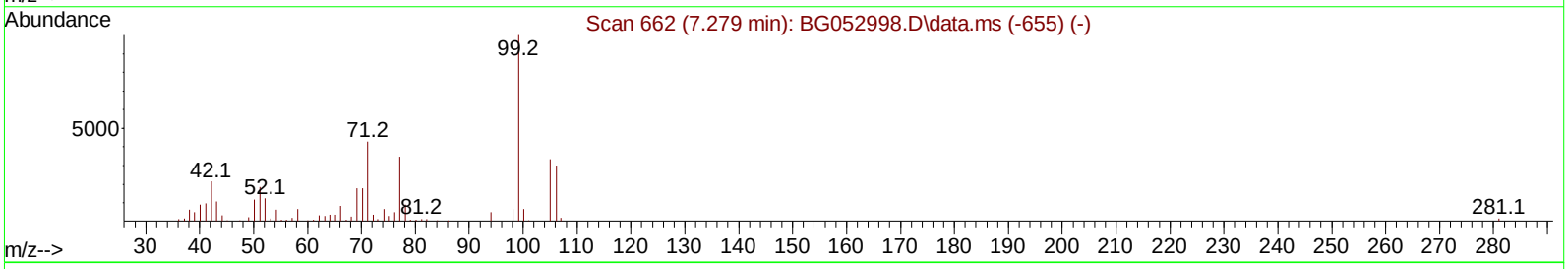

TIC: BG053000.D\data.ms

(7) Phenol-d5 (S)

7.279min (+ 0.002) 7.50 ng/ul m

response 19371

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	36.40	37.64
42.00	19.00	14.87#
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	8.137	152	28854	20.000	ng/ul	0.00
20) Naphthalene-d8	10.945	136	120757	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.751	164	92448	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.495	188	220051	20.000	ng/ul	0.00
79) Chrysene-d12	21.777	240	236720	20.000	ng/ul	# 0.00
88) Perylene-d12	25.084	264	225835	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.555	96	3344	4.094	ng/uL	0.00
4) Pyridine-d5	3.978	84	12898	6.325	ng/ul	0.01
7) Phenol-d5	7.279	99	19371m	7.496	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.449	67	55303	35.038	ng/ul	0.00
11) 2-Chlorophenol-d4	7.667	132	44693	25.469	ng/ul	0.00
15) 4-Methylphenol-d8	8.824	113	32286	17.246	ng/ul	0.00
21) Nitrobenzene-d5	9.294	128	32355	37.221	ng/ul	0.00
24) 2-Nitrophenol-d4	10.017	143	32968	35.958	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.557	165	57304	28.730	ng/ul	0.00
31) 4-Chloroaniline-d4	11.068	131	65801	24.961	ng/ul	0.00
46) Dimethylphthalate-d6	14.146	166	247049	34.785	ng/ul	0.00
49) Acenaphthylene-d8	14.446	160	273241	31.661	ng/ul	0.00
54) 4-Nitrophenol-d4	14.922	143	8698	7.284	ng/ul	0.00
60) Fluorene-d10	15.738	176	219874	35.258	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.844	200	45285	40.439	ng/ul	0.00
73) Anthracene-d10	17.595	188	398947	38.933	ng/ul	0.00
81) Pyrene-d10	19.874	212	496583	37.499	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.861	264	468094	40.114	ng/ul	0.00
Target Compounds						Qvalue

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

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