

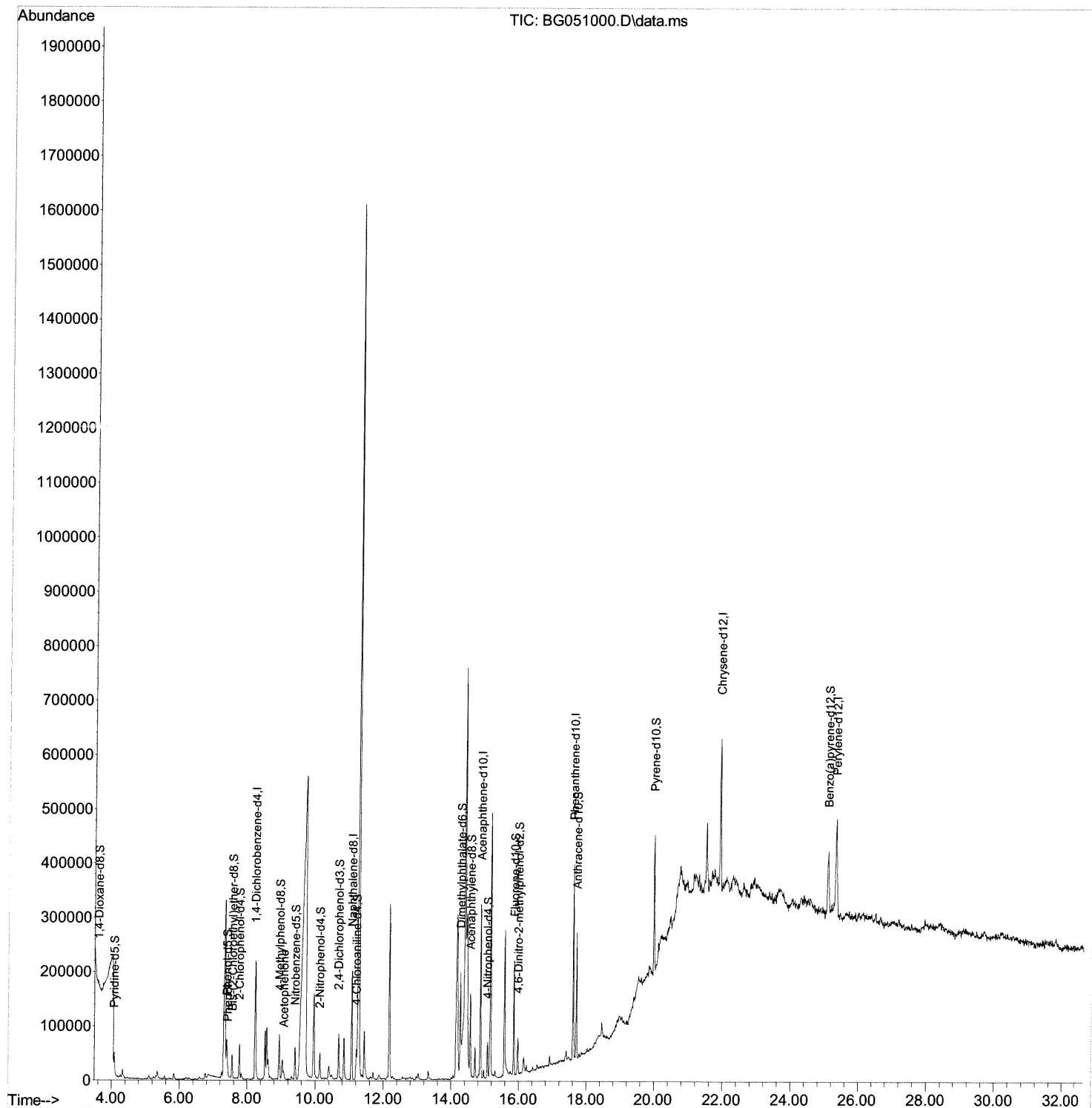
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG051000.D
Acq On : 12 Nov 2021 14:24
Operator : CG/JU
Sample : M4615-08
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0V14

Manual IntegrationsAPPROVED

Quant Time: Nov 12 14:58:29 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Nov 11 12:40:48 2021
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/12/2021
Supervised By :mohammad ahmed 11/17/2021



Quantitation Report (Qedit)

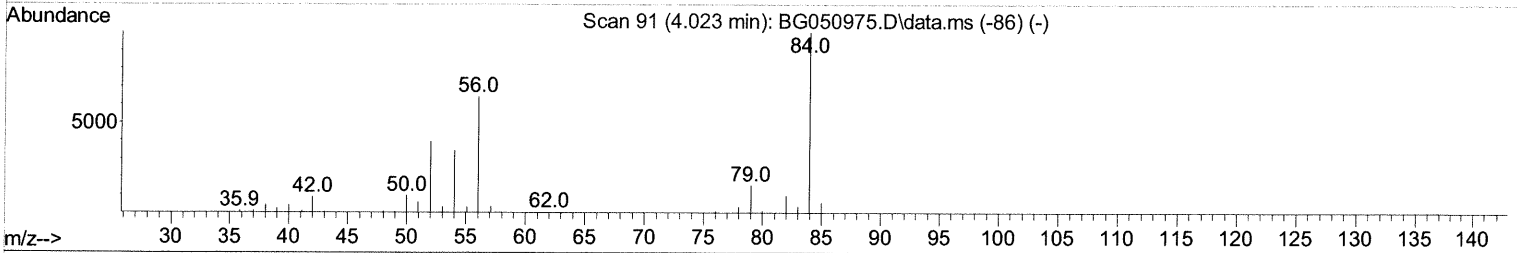
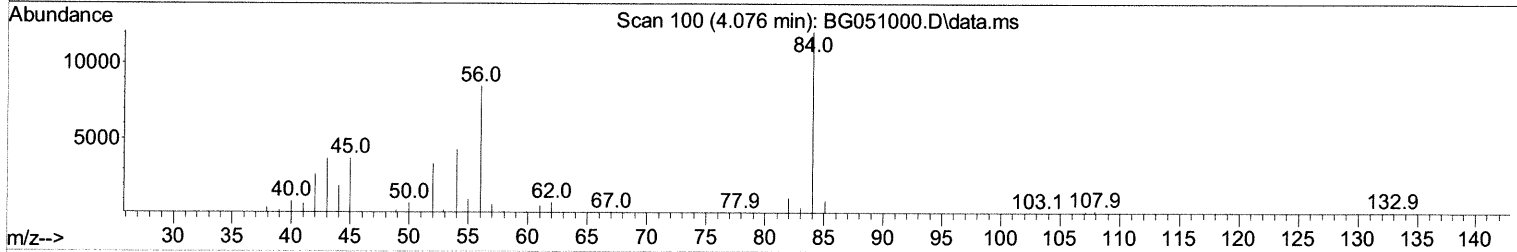
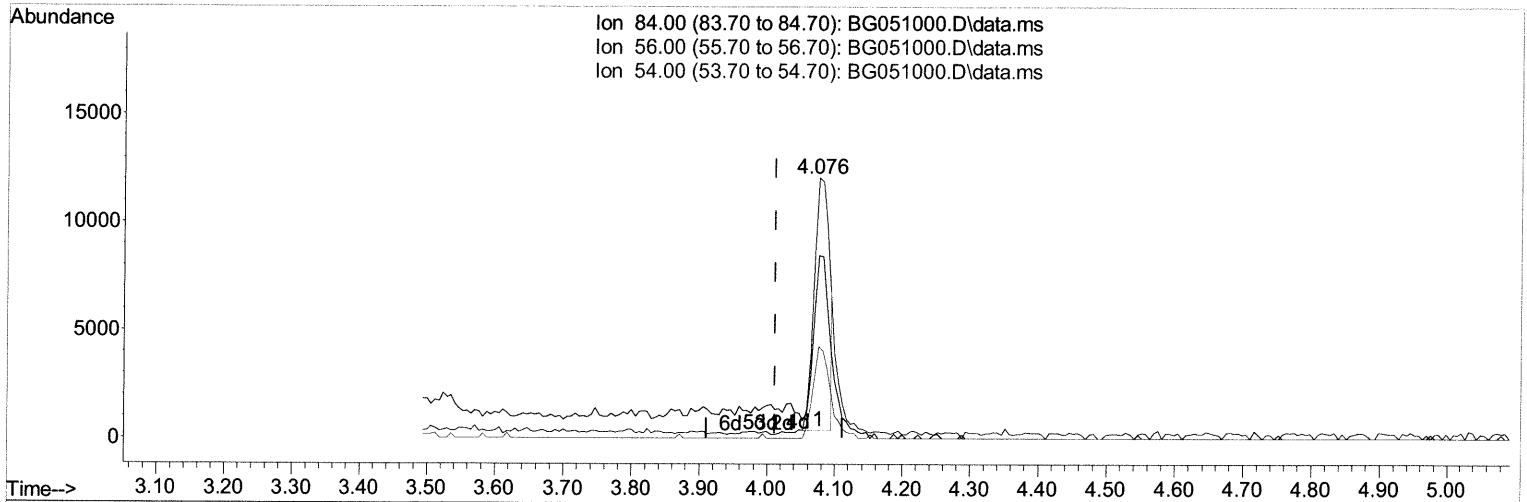
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TIC: BG051000.D\data.ms

(4) Pyridine-d5 (S)

4.076min (+ 0.065) 5.33 ng/ul

response 18419

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	70.12
54.00	31.50	35.33
0.00	0.00	0.00

Quantitation Report (Qedit)

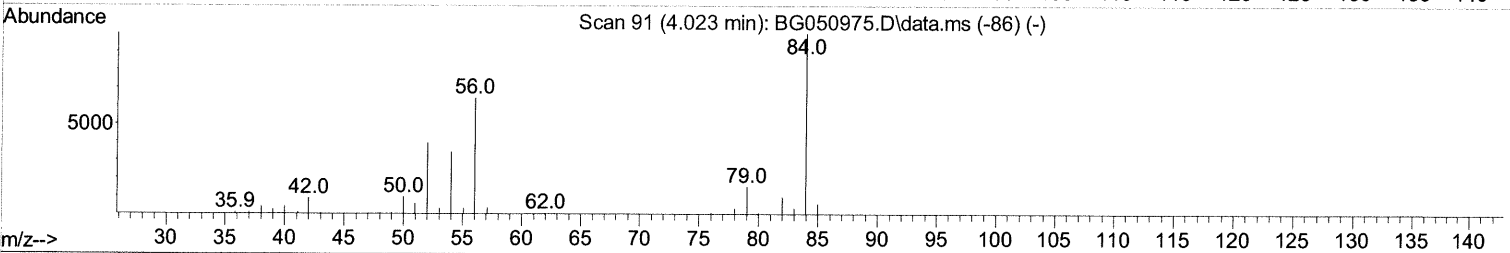
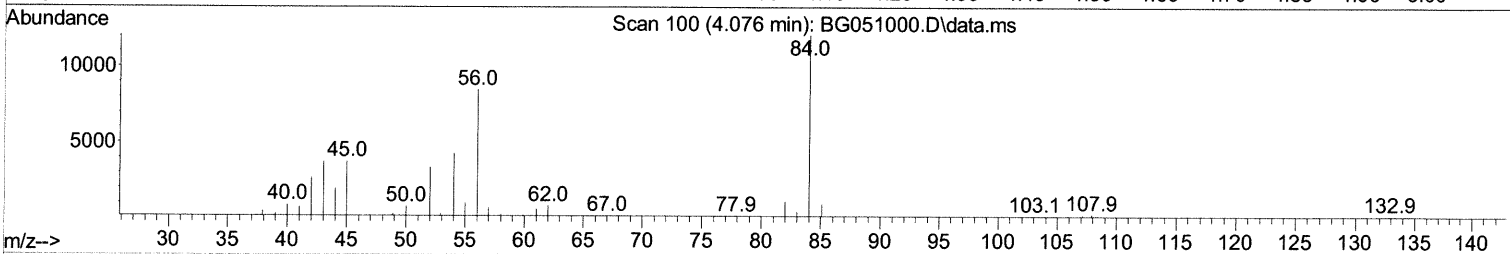
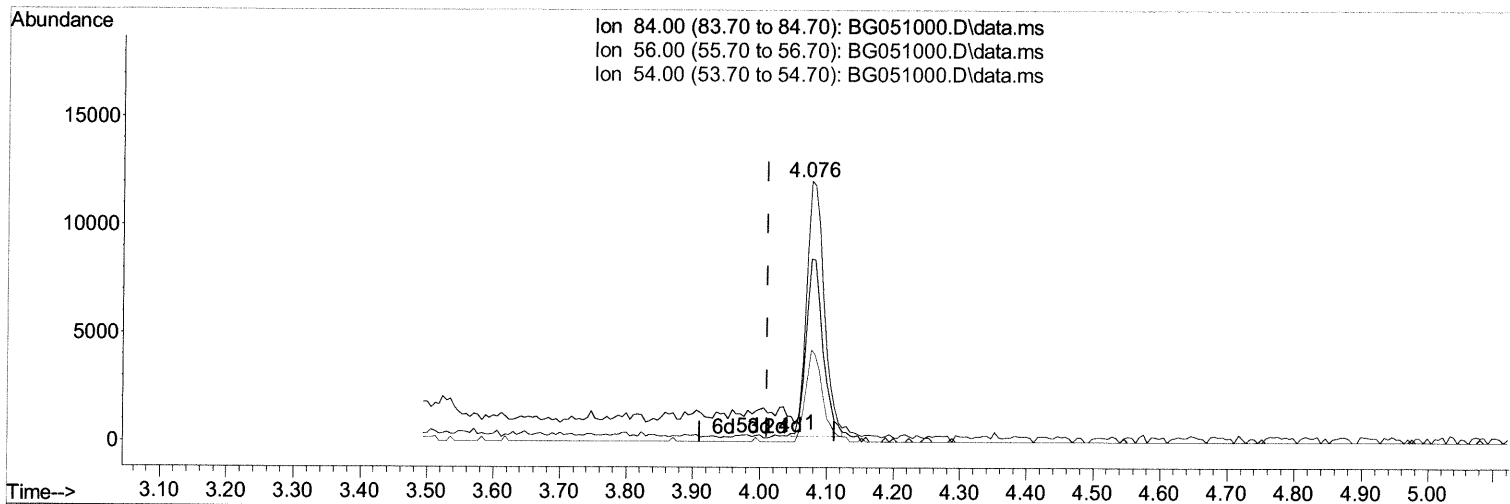
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TIC: BG051000.D\data.ms

(4) Pyridine-d5 (S)

4.076min (+ 0.065) 6.30 ng/ul m 11/17/21 Ju

response 21772

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	70.12
54.00	31.50	35.33
0.00	0.00	0.00

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 BNA_G
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/12/2021
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Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.242	152	37262	20.000 ng/ul	0.02
20) Naphthalene-d8	11.068	136	169134	20.000 ng/ul	0.02
38) Acenaphthene-d10	14.869	164	112859	20.000 ng/ul	0.02
64) Phenanthrene-d10	17.613	188	220597	20.000 ng/ul	0.02
79) Chrysene-d12	21.914	240	171959	20.000 ng/ul	0.02
88) Perylene-d12	25.333	264	179242	20.000 ng/ul	0.05
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.612	96	1563	1.354 ng/uL	0.02
4) Pyridine-d5	4.076	84	21772m	6.304 ng/ul	0.06
7) Phenol-d5	7.396	99	40619	10.218 ng/ul	0.03
9) Bis-(2-Chloroethyl)eth...	7.554	67	21010	8.182 ng/ul	0.01
11) 2-Chlorophenol-d4	7.772	132	28790	10.450 ng/ul	0.02
15) 4-Methylphenol-d8	8.953	113	34432	11.002 ng/ul	0.03
21) Nitrobenzene-d5	9.411	128	13998	9.739 ng/ul	0.01
24) 2-Nitrophenol-d4	10.139	143	15218	9.522 ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.686	165	31742	11.791 ng/ul	0.02
31) 4-Chloroaniline-d4	11.203	131	32860	8.060 ng/ul	0.02
46) Dimethylphthalate-d6	14.264	166	99329	11.504 ng/ul	0.02
49) Acenaphthylene-d8	14.564	160	110190	10.243 ng/ul	0.02
54) 4-Nitrophenol-d4	15.075	143	16349	10.443 ng/ul	0.04
60) Fluorene-d10	15.856	176	84122	10.998 ng/ul	0.02
65) 4,6-Dinitro-2-methylph...	15.974	200	13925	10.412 ng/ul	0.02
73) Anthracene-d10	17.713	188	137500	13.184 ng/ul	0.02
81) Pyrene-d10	19.987	212	142367	12.818 ng/ul	0.02
92) Benzo(a)pyrene-d12	25.092	264	121149	12.226 ng/ul	0.04
Target Compounds					
8) Phenol	7.425	94	6114	1.487 ng/ul#	87
16) Acetophenone	9.076	105	6620	1.362 ng/ul#	95

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