

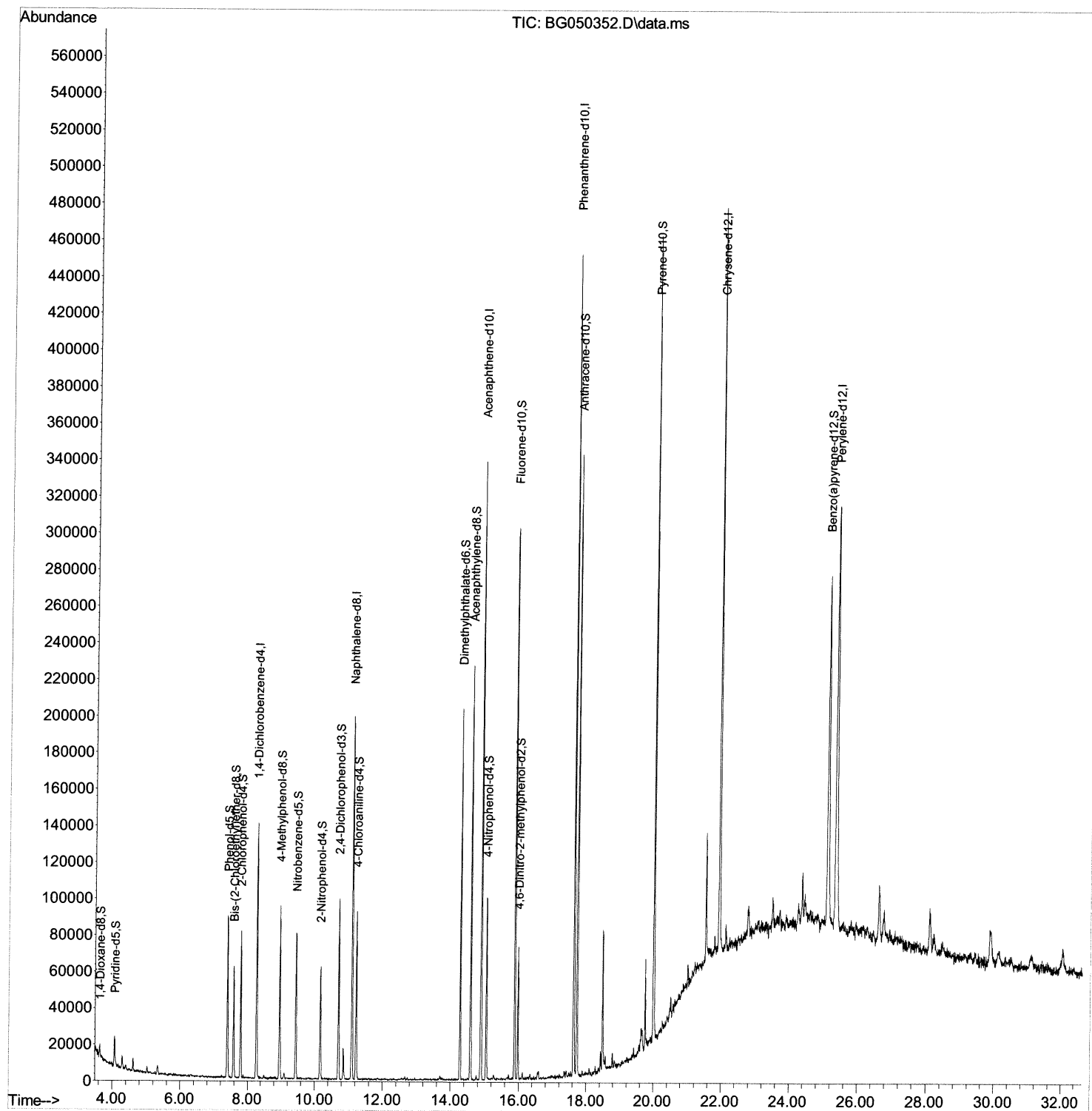
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
Data File : BG050352.D
Acq On : 1 Oct 2021 10:55
Operator : CG/JU
Sample : M3876-01
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7W1

Manual Integrations
APPROVED

mohammad
10/4/2021 11:50:04 AM

Quant Time: Oct 01 11:49:23 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Sep 30 16:27:24 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

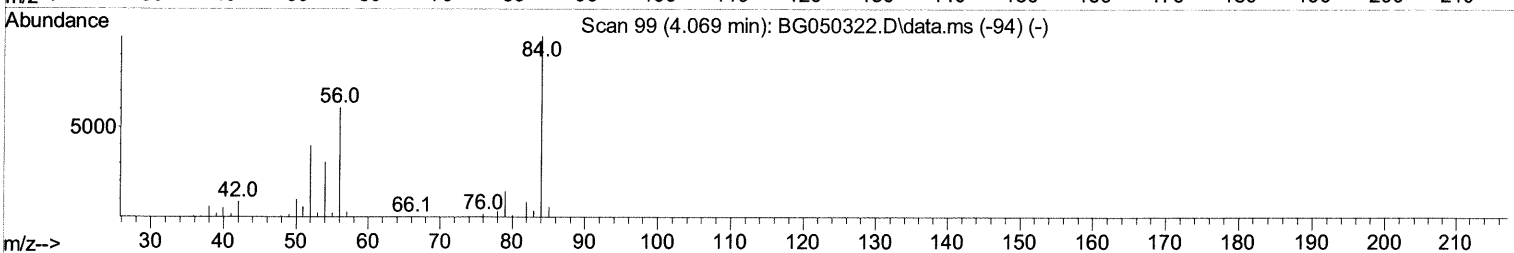
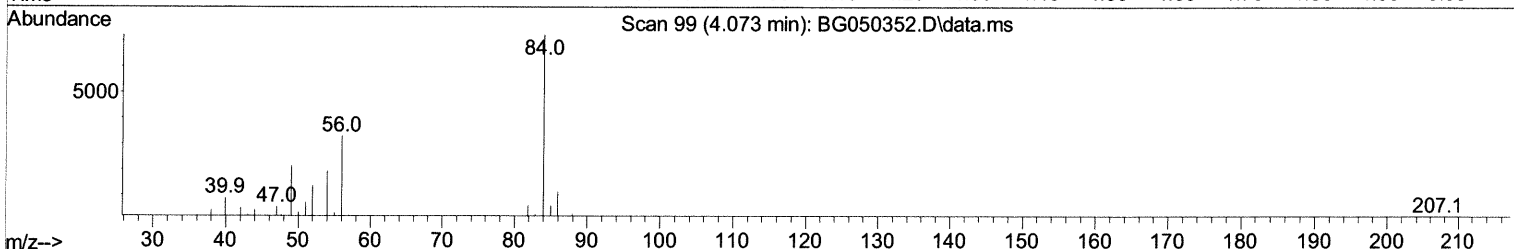
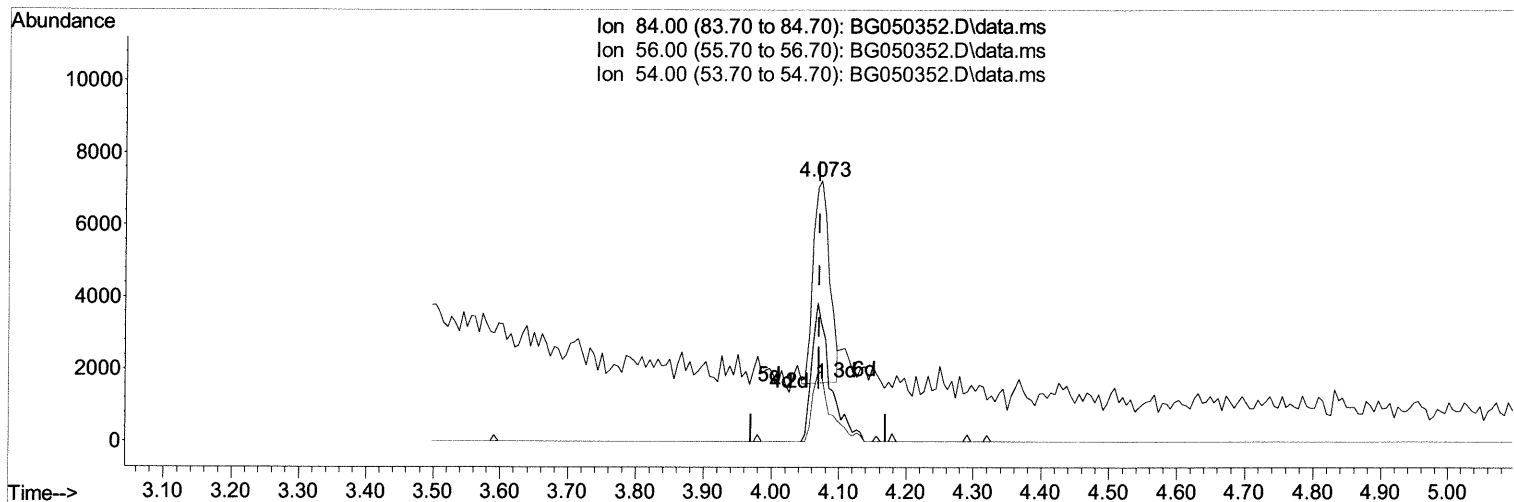
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TIC: BG050352.D\data.ms

(4) Pyridine-d5 (S)

4.073min (+ 0.004) 3.05 ng/ul

response 9483

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.00	45.42#
54.00	30.10	26.76
0.00	0.00	0.00

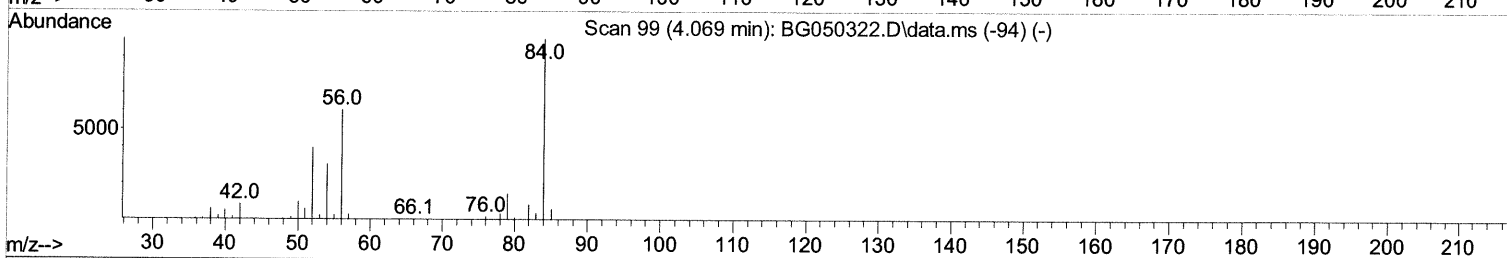
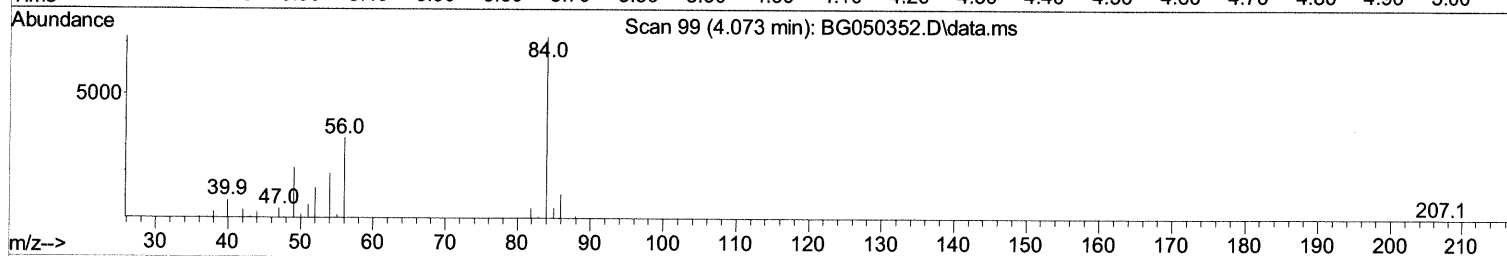
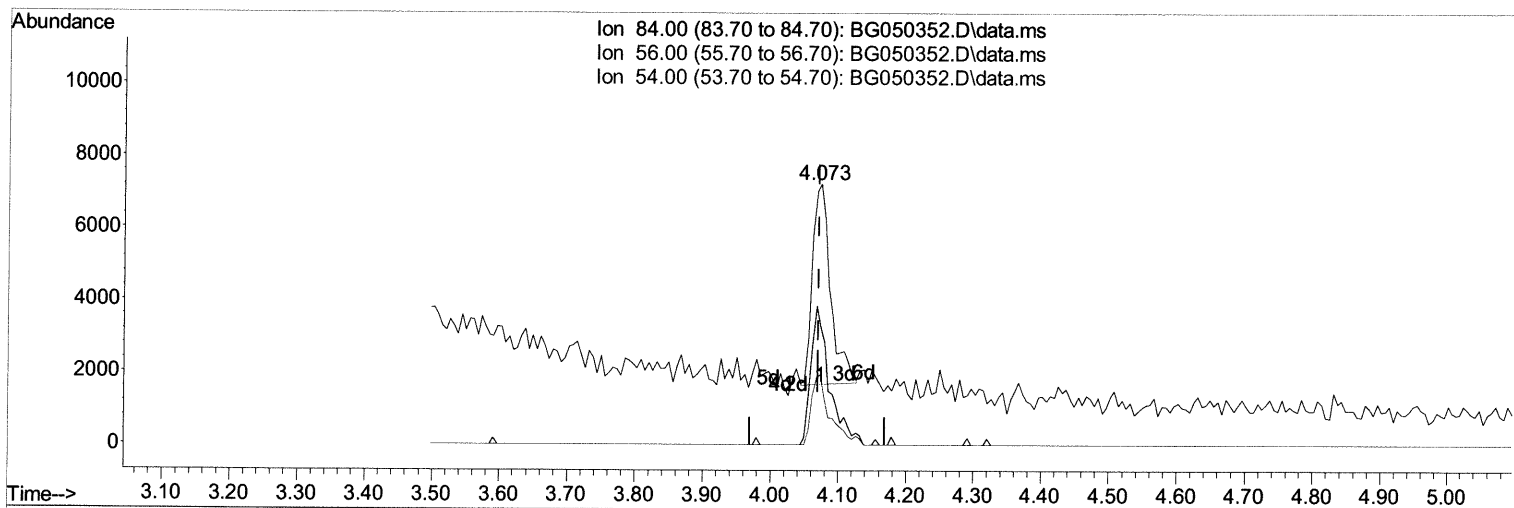
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TIC: BG050352.D\data.ms

(4) Pyridine-d5 (S)

4.073min (+ 0.004) 3.29 ng/ul m 10/05/21 JU

response 10236

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.00	45.42#
54.00	30.10	26.76
0.00	0.00	0.00

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Instrument :
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Manual Integrations
 APPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.268	152	38381	20.000	ng/ul	0.00
20) Naphthalene-d8	11.100	136	162856	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.890	164	117229	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.628	188	275879	20.000	ng/ul	0.00
79) Chrysene-d12	21.928	240	252513	20.000	ng/ul	-0.02
88) Perylene-d12	25.360	264	246515	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.638	96	2695	2.580	ng/uL	0.00
4) Pyridine-d5	4.073	84	10236m	3.288	ng/ul	> 0.00
7) Phenol-d5	7.398	99	48382	14.428	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.586	67	32519	15.409	ng/ul	0.00
11) 2-Chlorophenol-d4	7.792	132	33930	14.441	ng/ul	0.00
15) 4-Methylphenol-d8	8.955	113	35416	13.954	ng/ul	0.00
21) Nitrobenzene-d5	9.443	128	18915	16.603	ng/ul	0.00
24) 2-Nitrophenol-d4	10.166	143	19704	16.115	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.700	165	34218	14.172	ng/ul	0.00
31) 4-Chloroaniline-d4	11.223	131	47768	13.849	ng/ul	0.00
46) Dimethylphthalate-d6	14.279	166	126723	15.795	ng/ul	-0.01
49) Acenaphthylene-d8	14.584	160	162376	16.116	ng/ul	0.00
54) 4-Nitrophenol-d4	15.048	143	22515	14.571	ng/ul	0.00
60) Fluorene-d10	15.877	176	121697	16.836	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.977	200	17492	12.712	ng/ul	0.00
73) Anthracene-d10	17.728	188	204475	17.544	ng/ul	0.00
81) Pyrene-d10	20.001	212	242896	18.090	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.119	264	201093	16.223	ng/ul	-0.02

Target Compounds Qvalue

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