

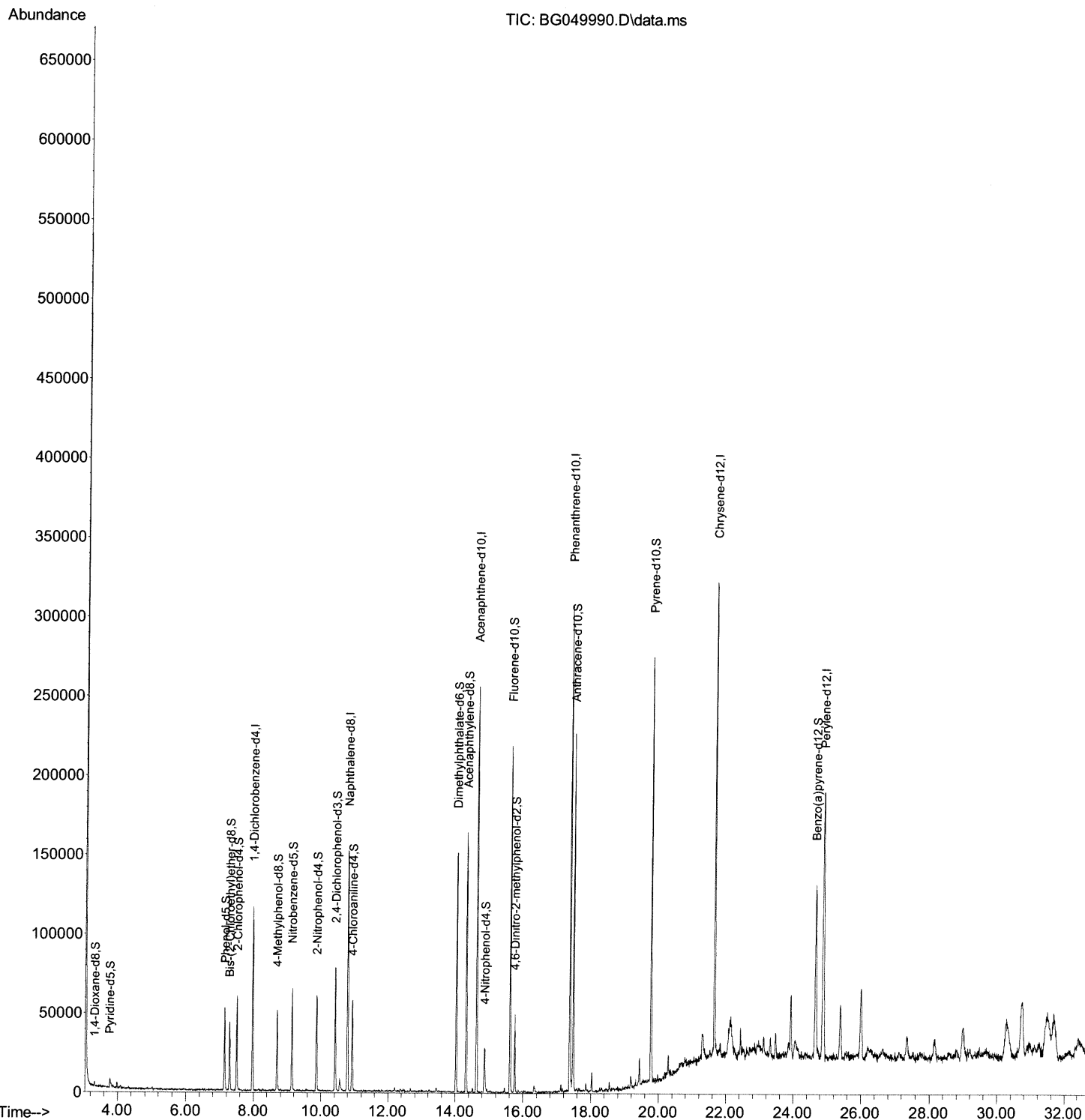
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082621\
Data File : BG049990.D
Acq On : 27 Aug 2021 1:04
Operator : CG/JU
Sample : M3458-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AR4

Manual Integrations
APPROVED

mohammad
8/27/2021 5:09:01 PM

Quant Time: Aug 27 02:02:32 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 26 15:41:36 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

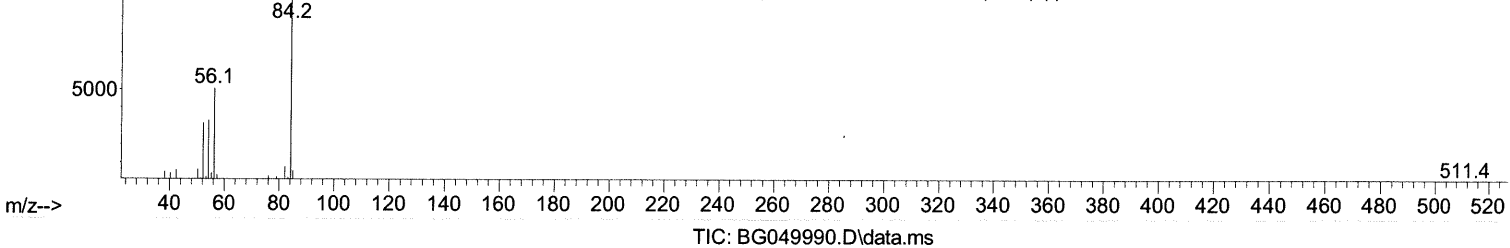
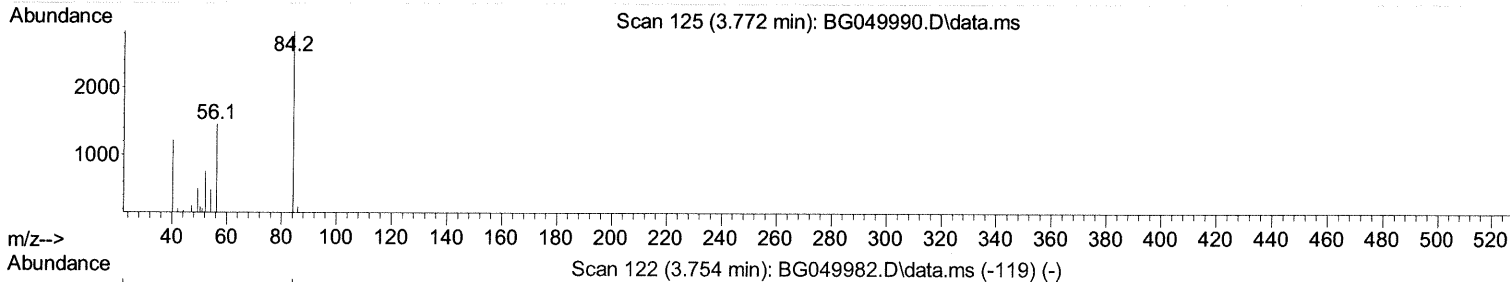
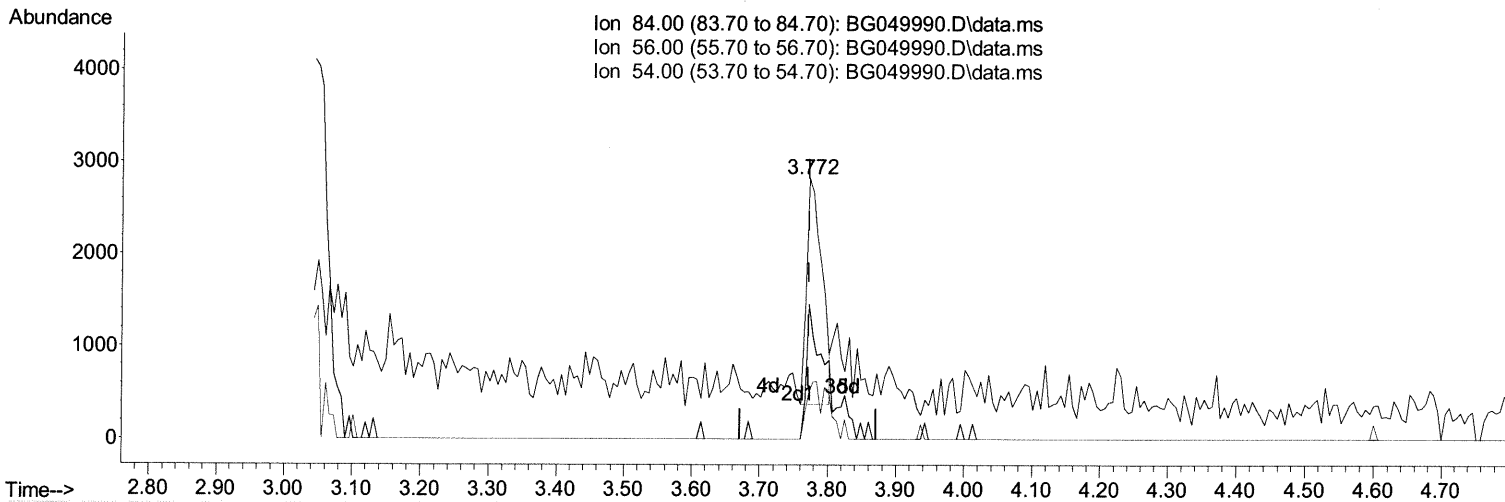
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(4) Pyridine-d5 (S)

3.772min (+ 0.000) 2.10 ng/ul

response 3861

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	54.30	51.70
54.00	37.00	17.30#
0.00	0.00	0.00

Quantitation Report (Qedit)

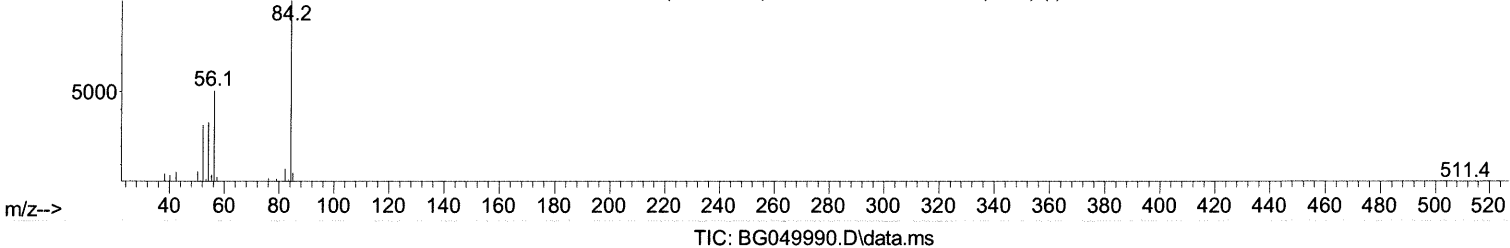
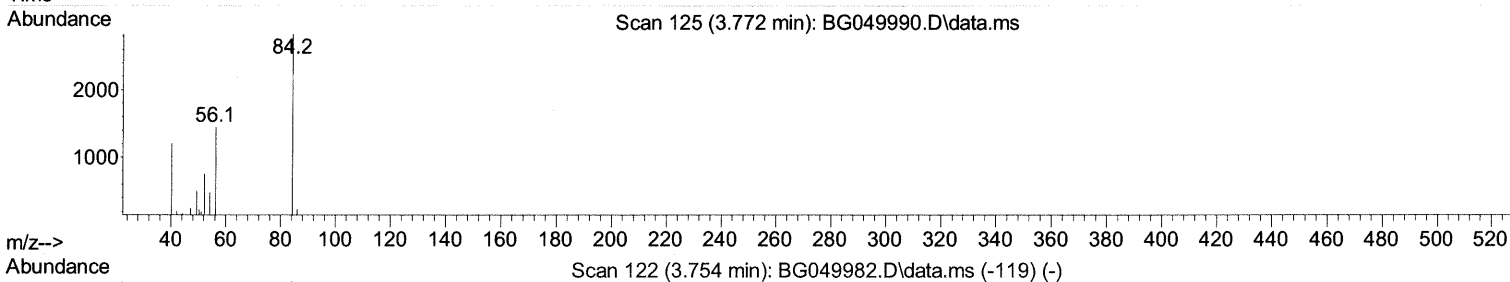
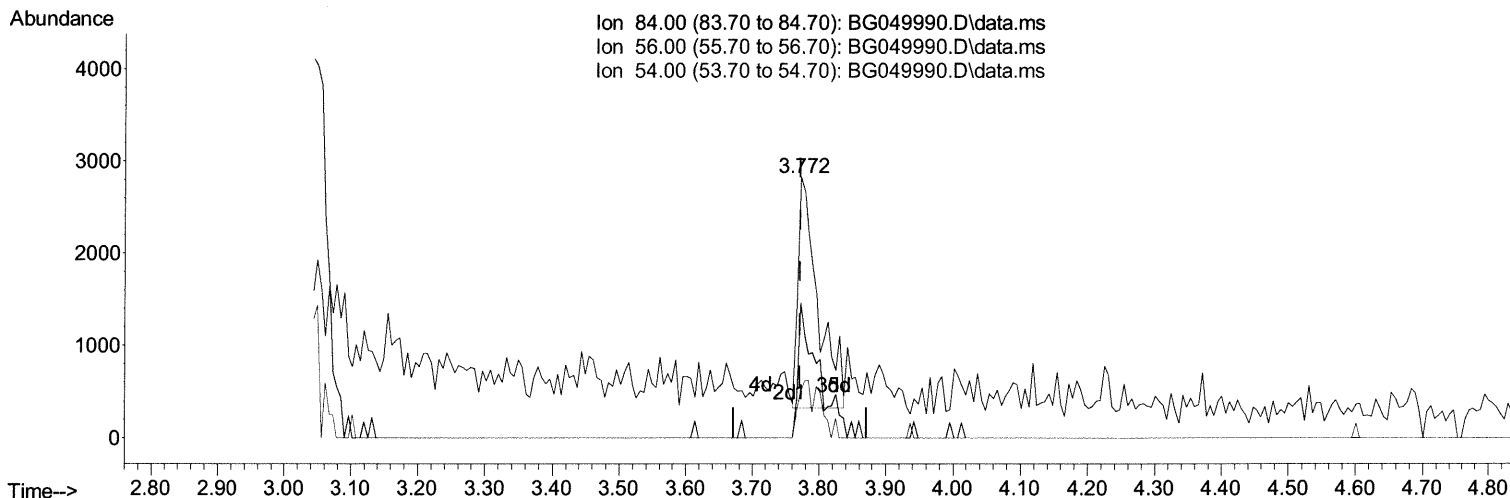
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(4) Pyridine-d5 (S)

3.772min (+ 0.000) 2.84 ng/ul m 08/30/21 ju

response 5226

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	54.30	51.70
54.00	37.00	17.30#
0.00	0.00	0.00

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.966	152	31503	20.000 ng/ul	-0.02
20) Naphthalene-d8	10.786	136	124488	20.000 ng/ul	-0.02
38) Acenaphthene-d10	14.628	164	88097	20.000 ng/ul	-0.01
64) Phenanthrene-d10	17.383	188	186125	20.000 ng/ul	-0.01
79) Chrysene-d12	21.653	240	176376	20.000 ng/ul	-0.01
88) Perylene-d12	24.884	264	176356	20.000 ng/ul	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.325	96	1239	2.028 ng/ul	-0.02
4) Pyridine-d5	3.772	84	5226m	2.840 ng/ul	> 0.00
7) Phenol-d5	7.138	99	30053	11.234 ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.285	67	20738	13.954 ng/ul	-0.02
11) 2-Chlorophenol-d4	7.502	132	28071	14.024 ng/ul	-0.01
15) 4-Methylphenol-d8	8.695	113	20944	9.941 ng/ul	0.00
21) Nitrobenzene-d5	9.135	128	14474	14.997 ng/ul	-0.02
24) 2-Nitrophenol-d4	9.869	143	17628	15.339 ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.422	165	29342	14.315 ng/ul	0.00
31) 4-Chloroaniline-d4	10.927	131	30349	10.878 ng/ul	-0.01
46) Dimethylphthalate-d6	14.029	166	100025	14.836 ng/ul	-0.01
49) Acenaphthylene-d8	14.322	160	121933	15.424 ng/ul	-0.01
54) 4-Nitrophenol-d4	14.869	143	9054	9.282 ng/ul	0.01
60) Fluorene-d10	15.620	176	87077	15.233 ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.756	200	11500	9.604 ng/ul	-0.01
73) Anthracene-d10	17.477	188	132769	15.885 ng/ul	-0.02
81) Pyrene-d10	19.768	212	150980	15.841 ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.655	264	115237	12.317 ng/ul	-0.02

Target Compounds Qvalue

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