

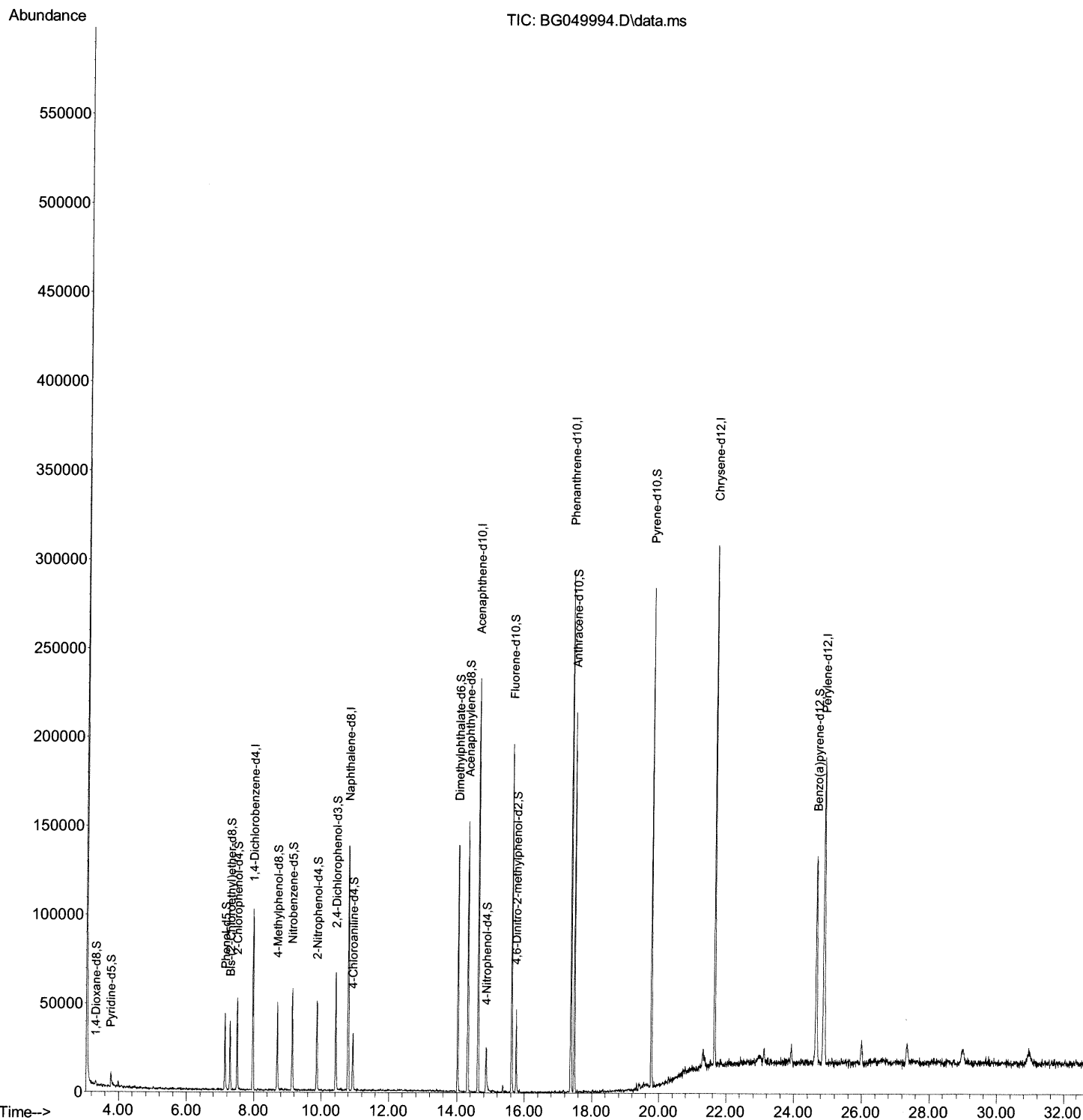
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082621\
Data File : BG049994.D
Acq On : 27 Aug 2021 3:49
Operator : CG/JU
Sample : M3458-03
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AR2

Manual Integrations
APPROVED

mohammad
8/27/2021 5:08:57 PM

Quant Time: Aug 27 04:30:44 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)

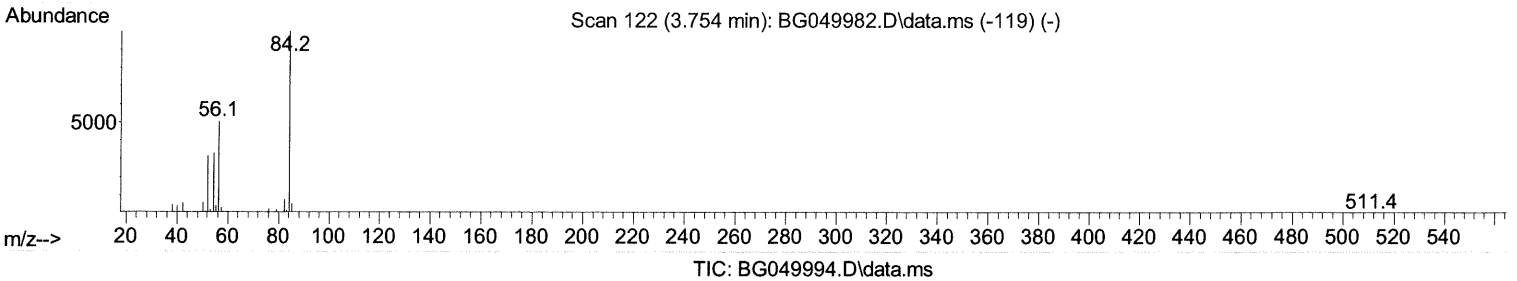
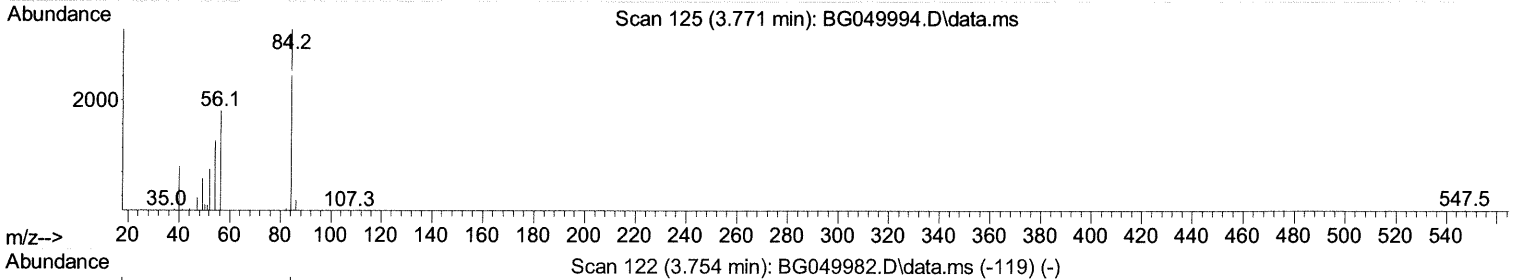
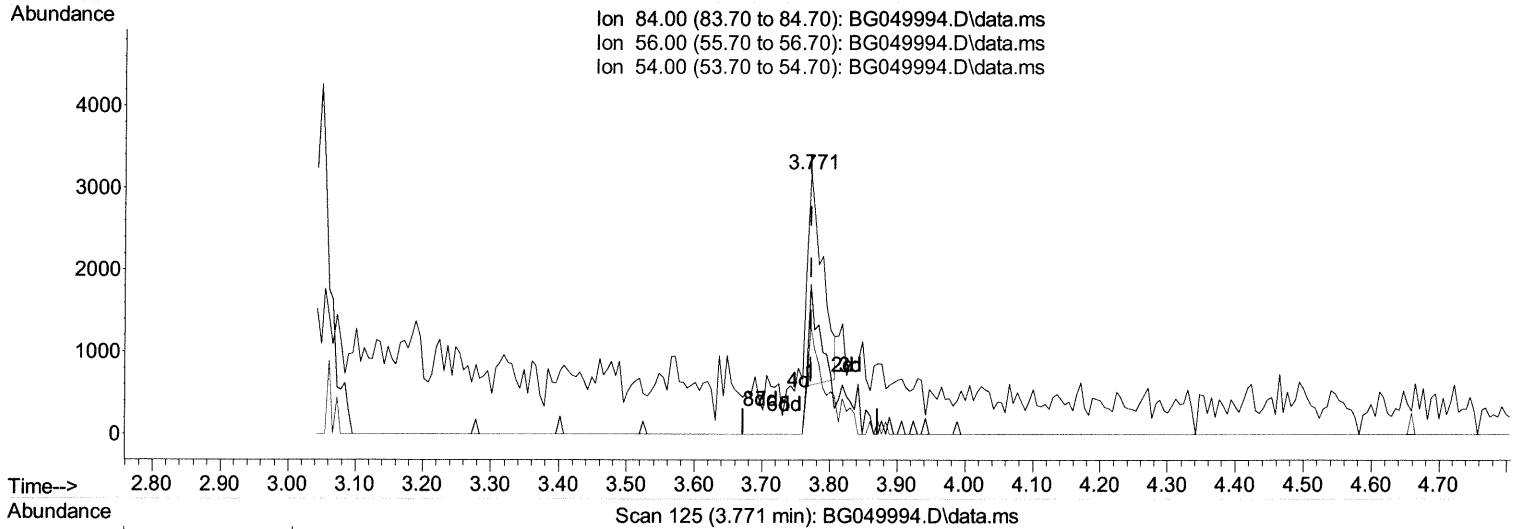
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(4) Pyridine-d5 (S)

3.771min (-0.000) 2.38 ng/ul

response 3951

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	54.30	57.37
54.00	37.00	41.50
0.00	0.00	0.00

Quantitation Report (Qedit)

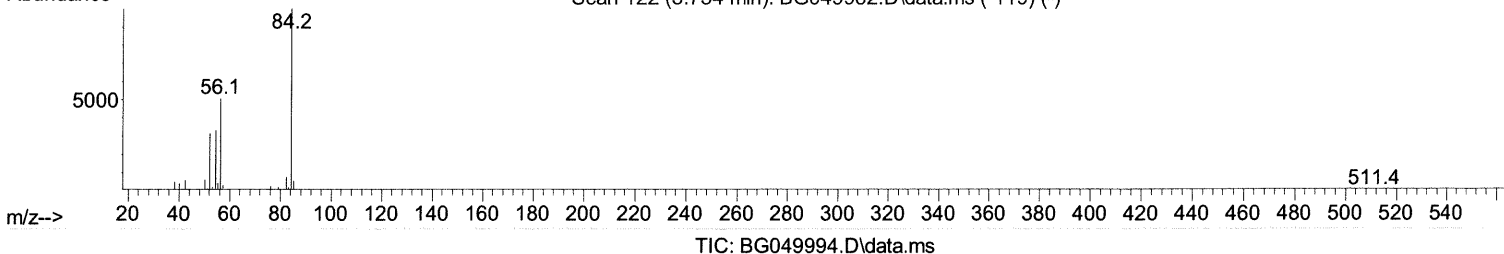
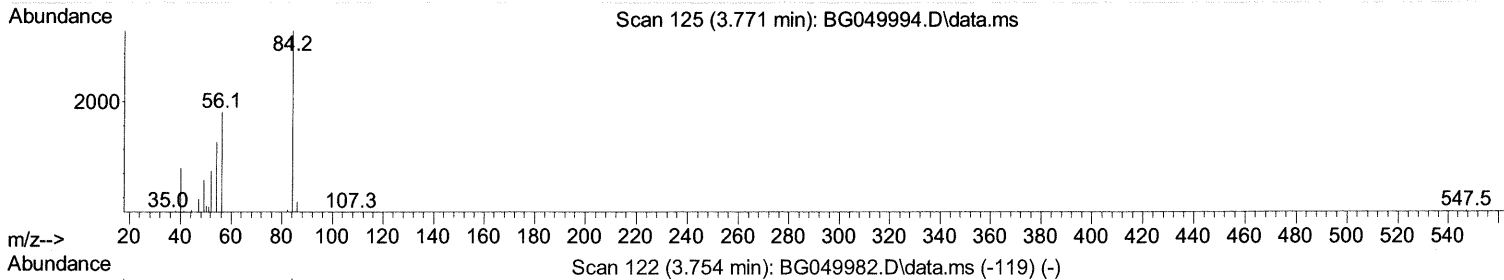
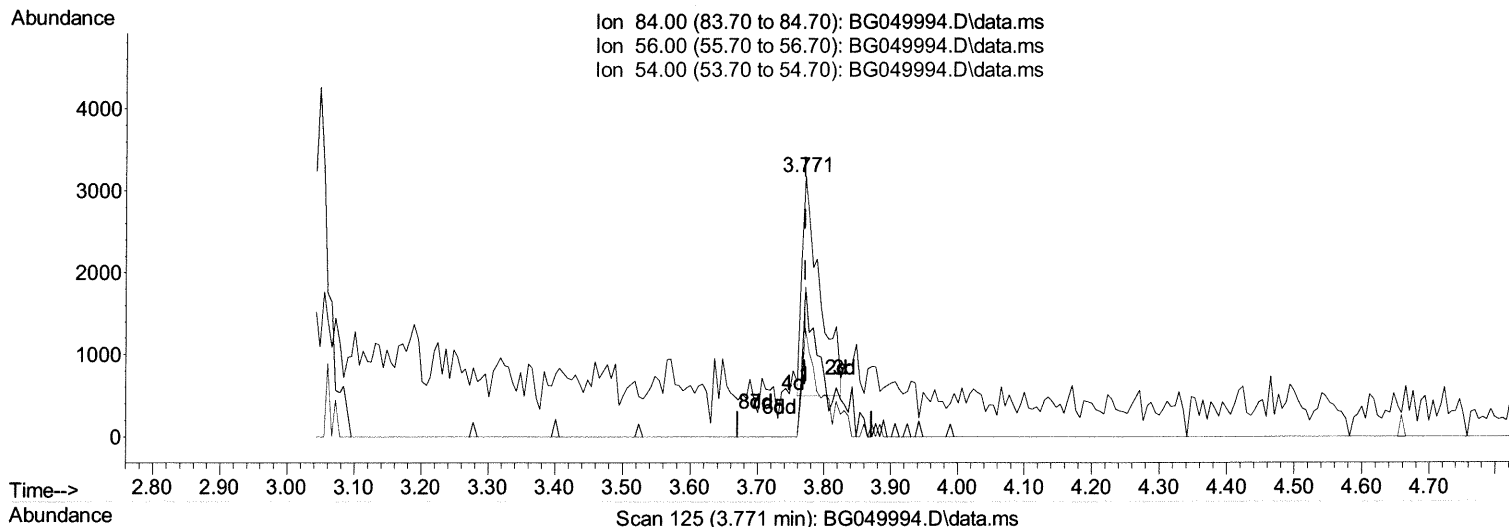
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(4) Pyridine-d5 (S)

3.771min (-0.000) 2.97 ng/ul m 08/30/21JU

response 4923

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	54.30	57.37
54.00	37.00	41.50
0.00	0.00	0.00

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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	7.971	152	28380	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.785	136	117079	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.627	164	82434	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.382	188	183567	20.000	ng/ul	-0.01
79) Chrysene-d12	21.653	240	178251	20.000	ng/ul	-0.01
88) Perylene-d12	24.878	264	175365	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.325	96	1313	2.386	ng/ul	-0.02
4) Pyridine-d5	3.771	84	4923m >	2.970	ng/ul >	0.00
7) Phenol-d5	7.137	99	26341	10.930	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.290	67	19105	14.270	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.496	132	24101	13.366	ng/ul	-0.02
15) 4-Methylphenol-d8	8.688	113	19010	10.016	ng/ul	-0.01
21) Nitrobenzene-d5	9.140	128	13881	15.293	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.869	143	14931	13.815	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.421	165	25066	13.003	ng/ul	0.00
31) 4-Chloroaniline-d4	10.926	131	17293	6.591	ng/ul	-0.01
46) Dimethylphthalate-d6	14.022	166	93727	14.857	ng/ul	-0.02
49) Acenaphthylene-d8	14.322	160	108967	14.731	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.868	143	6898	7.557	ng/ul	0.01
60) Fluorene-d10	15.620	176	77493	14.488	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.755	200	10786	9.133	ng/ul	-0.01
73) Anthracene-d10	17.482	188	129414	15.700	ng/ul	-0.01
81) Pyrene-d10	19.773	212	162109	16.829	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.661	264	126847	13.634	ng/ul	-0.01

Target Compounds Qvalue

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