

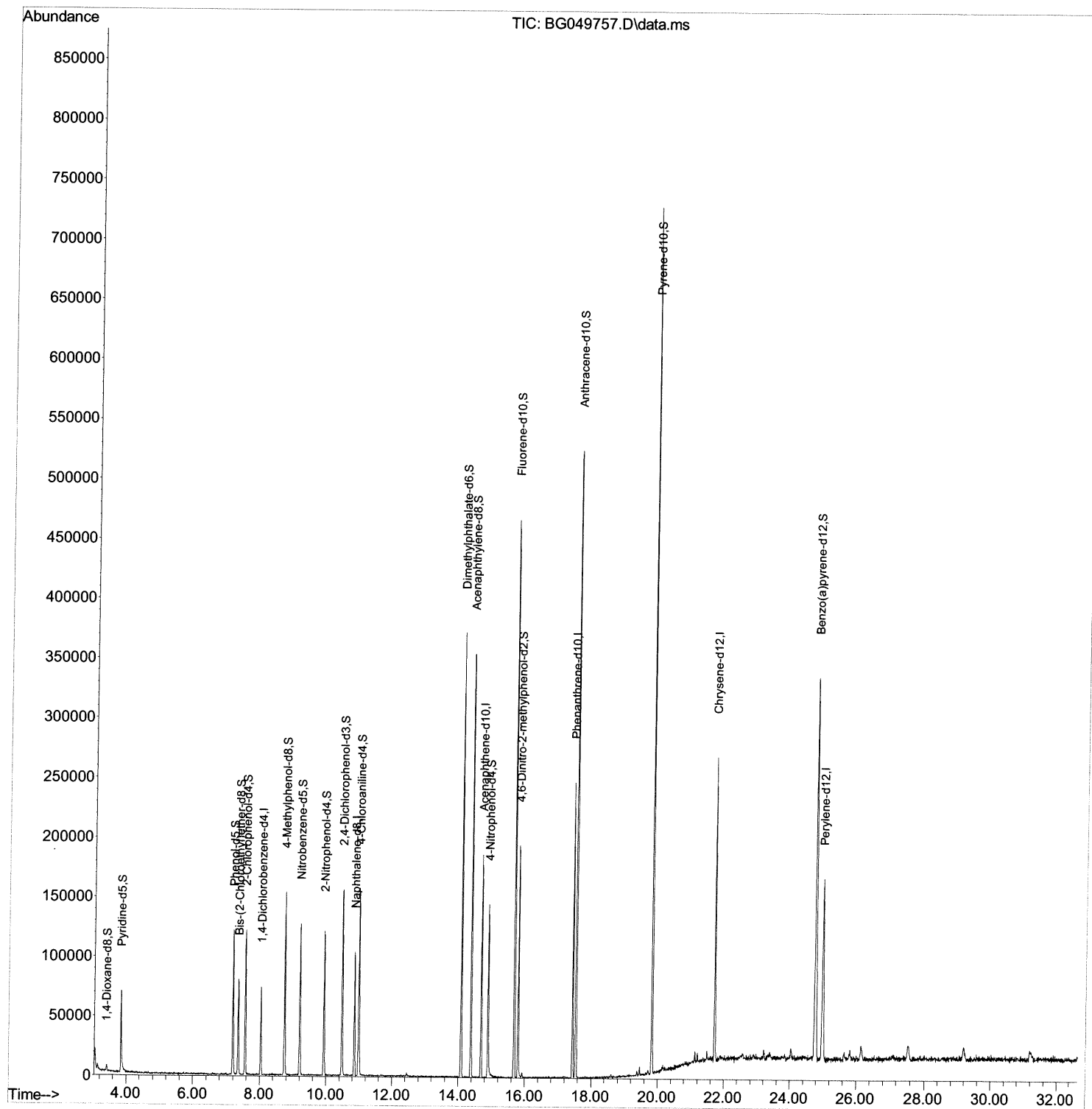
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081021\
Data File : BG049757.D
Acq On : 10 Aug 2021 15:00
Operator : CG/JU
Sample : PB138137BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK137

Manual Integrations
APPROVED

mohammad
8/11/2021 9:31:57 PM

Quant Time: Aug 10 17:07:14 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 10 16:55:58 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

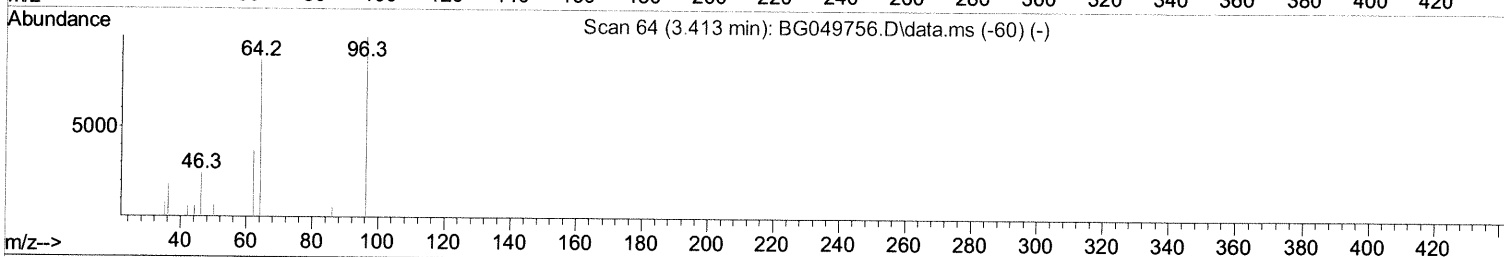
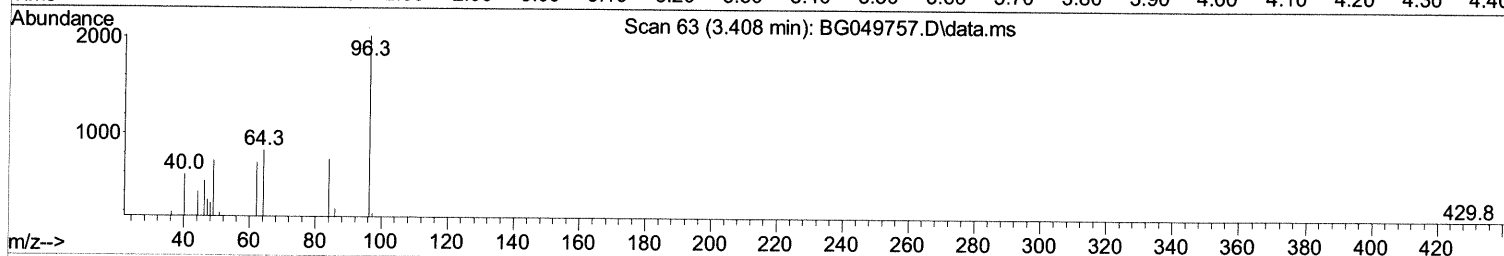
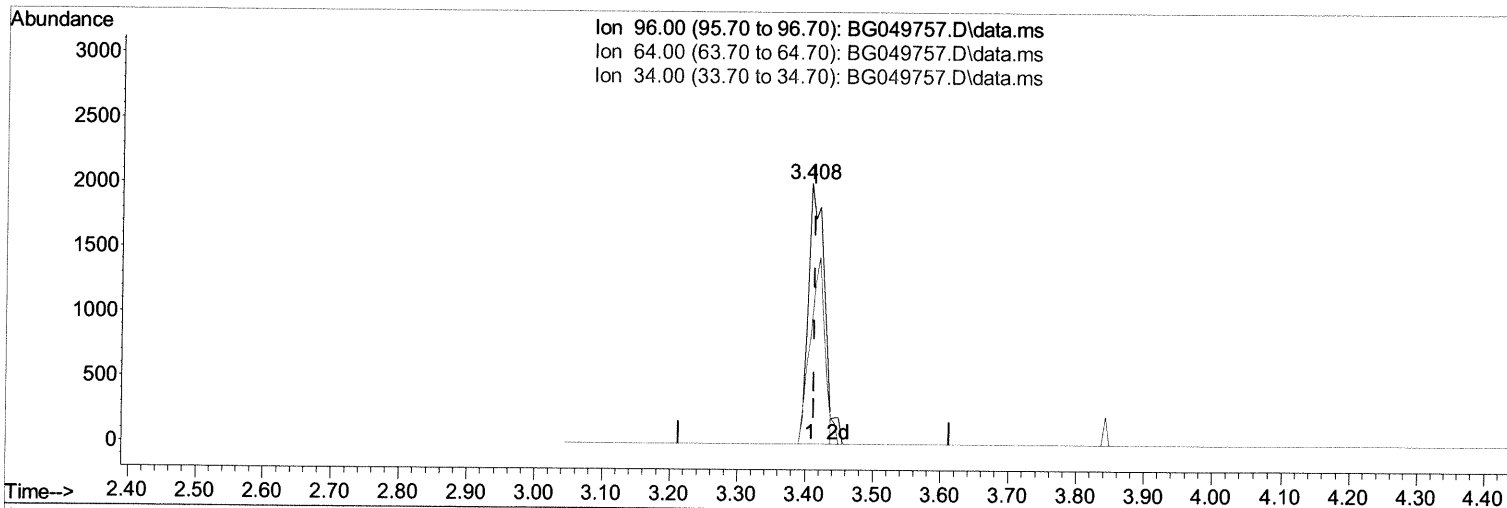
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TIC: BG049757.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.408min (-0.005) 7.49 ng/uL

response 3177

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.10	41.50#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

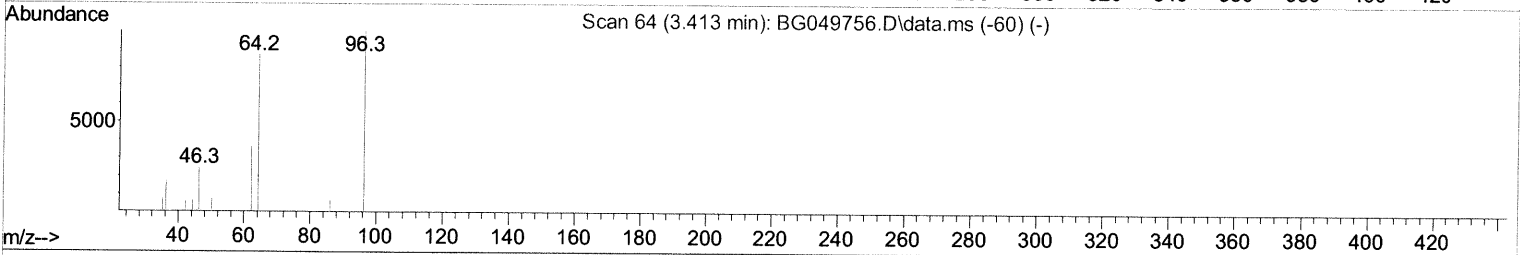
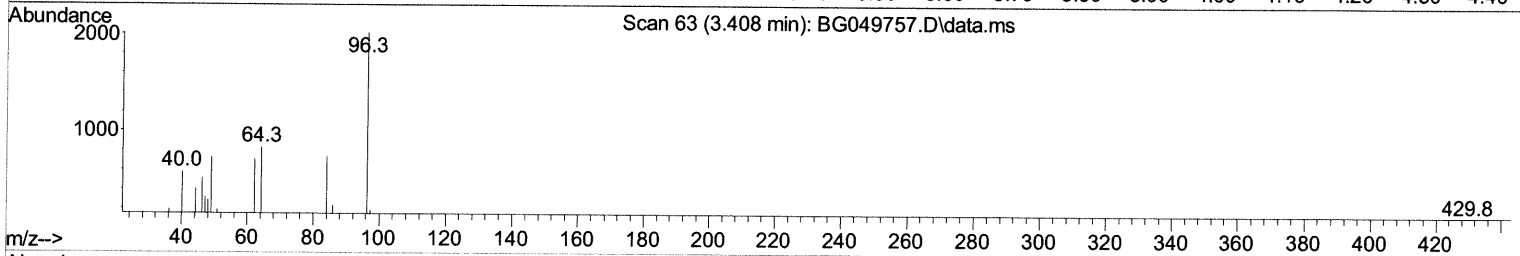
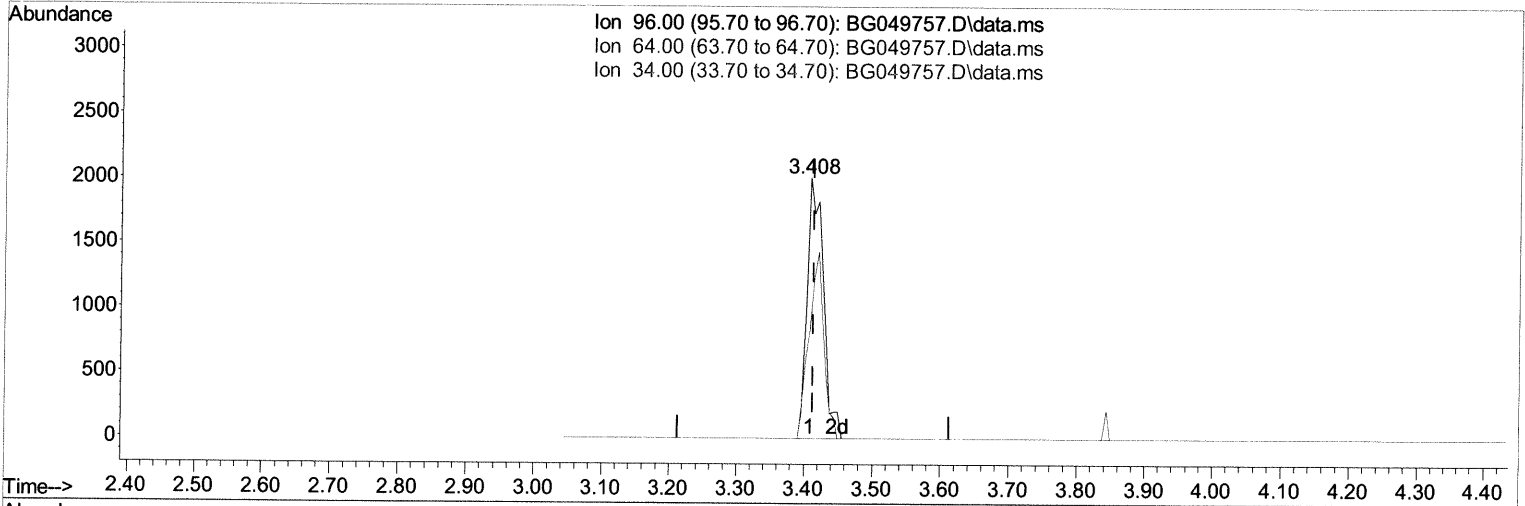
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Manual Integrations
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TIC: BG049757.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.408min (-0.005) 7.83 ng/uL m 08/12/21 JU

response 3318

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.10	41.50#
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081021\
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 Operator : CG/JU
 Sample : PB138137BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual Integrations
 APPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.037	152	19824	20.000	ng/ul	0.00
20) Naphthalene-d8	10.857	136	85814	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.687	164	66929	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.442	188	156577	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.718	240	155683	20.000	ng/ul	0.00
88) Perylene-d12	24.991	264	152396	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.408	96	3318m	7.825	ng/ul	0.0008112134
4) Pyridine-d5	3.837	84	44852	35.251	ng/ul	0.00
7) Phenol-d5	7.197	99	66346	36.876	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.356	67	38835	37.578	ng/ul	0.00
11) 2-Chlorophenol-d4	7.567	132	49973	38.808	ng/ul	0.00
15) 4-Methylphenol-d8	8.748	113	53904	39.295	ng/ul	0.00
21) Nitrobenzene-d5	9.206	128	28191	41.704	ng/ul	0.00
24) 2-Nitrophenol-d4	9.929	143	33817	43.805	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.481	165	56092	38.534	ng/ul	0.00
31) 4-Chloroaniline-d4	10.992	131	80370	40.331	ng/ul	0.00
46) Dimethylphthalate-d6	14.088	166	238584	45.425	ng/ul	0.00
49) Acenaphthylene-d8	14.381	160	254820	40.779	ng/ul	0.00
54) 4-Nitrophenol-d4	14.892	143	34436	40.551	ng/ul	0.00
60) Fluorene-d10	15.680	176	191617	42.442	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.803	200	42092	39.942	ng/ul	0.00
73) Anthracene-d10	17.542	188	317719	42.987	ng/ul	0.00
81) Pyrene-d10	19.827	212	386685	45.079	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.773	264	348027	41.678	ng/ul	0.00

Target Compounds Qvalue

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