

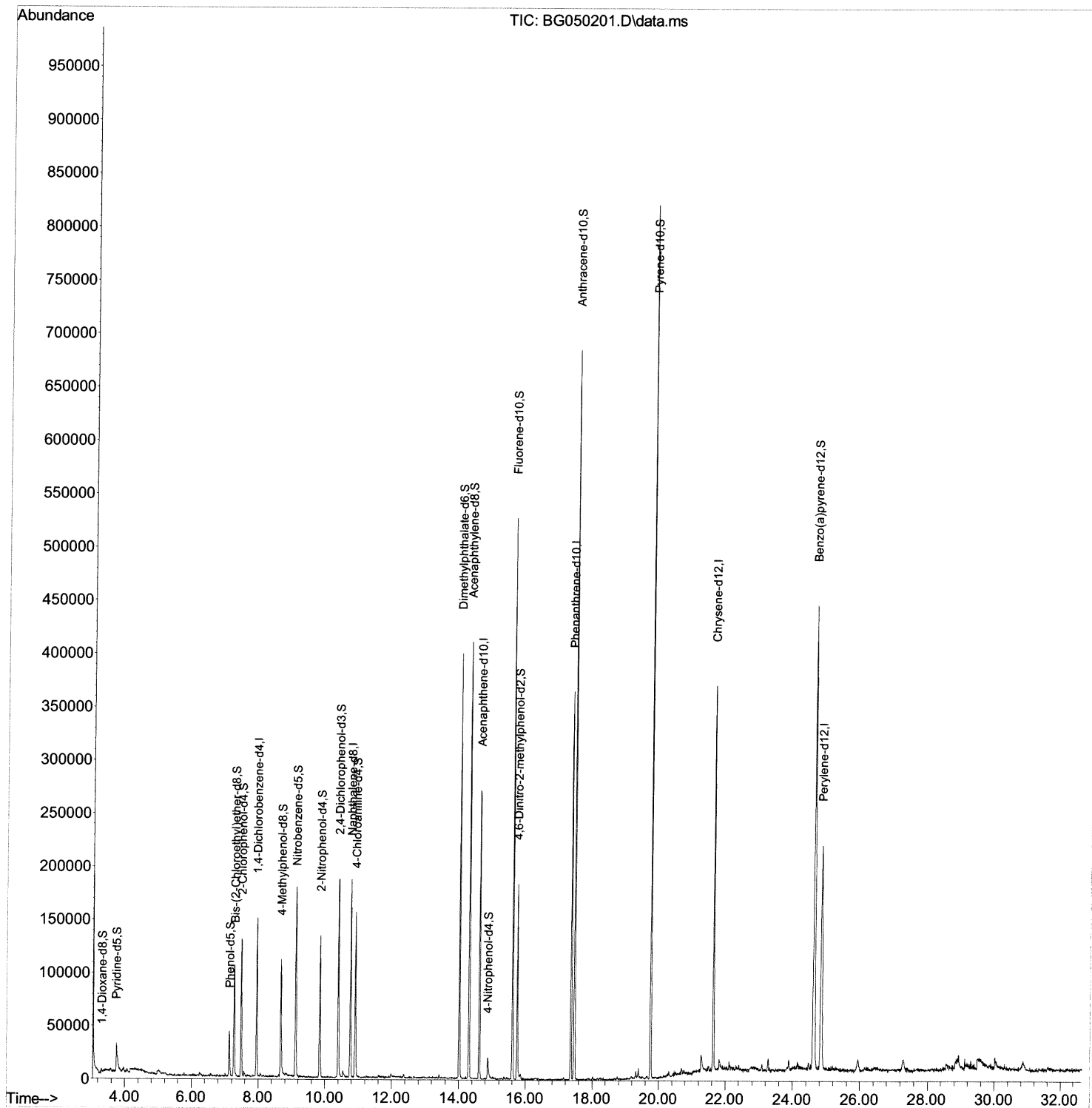
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050201.D
Acq On : 8 Sep 2021 15:54
Operator : CG/JU
Sample : M3608-15
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKL1

Manual Integrations
APPROVED

mohammad
9/9/2021 8:41:22 AM

Quant Time: Sep 08 16:50:43 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

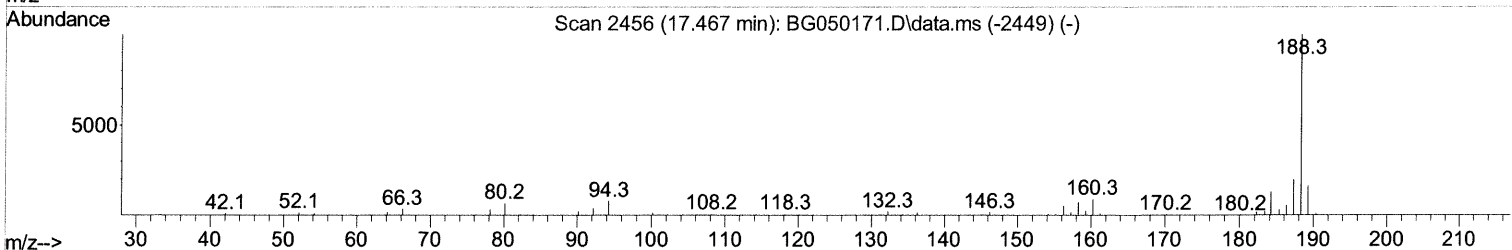
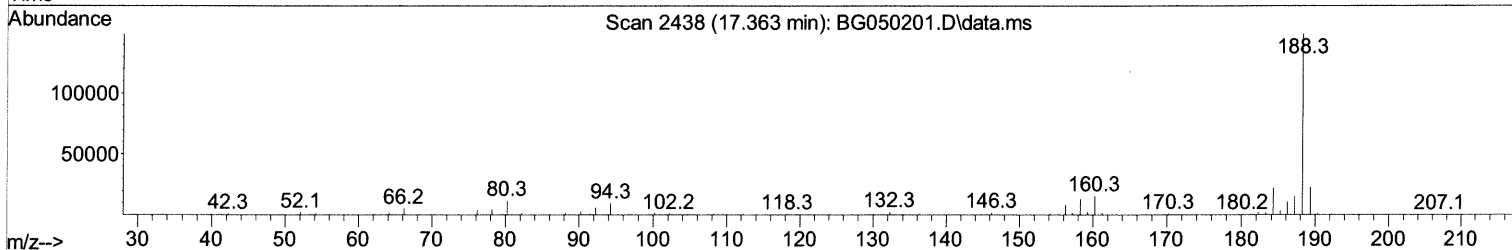
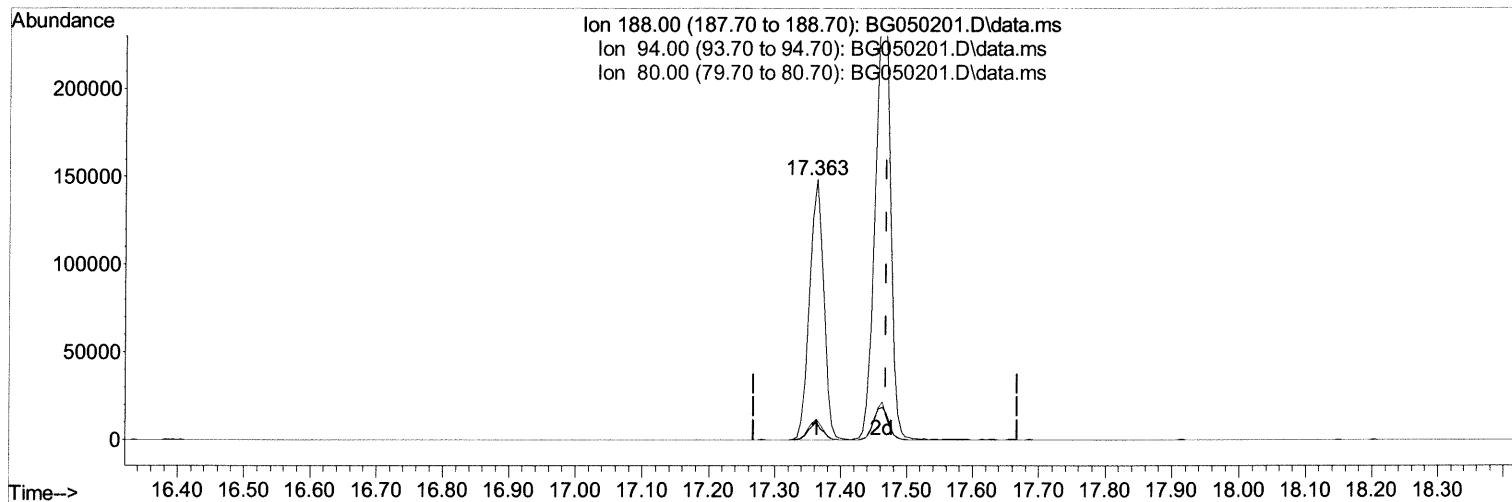
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TIC: BG050201.D\data.ms

(73) Anthracene-d10 (S)

17.363min (-0.105) 22.28 ng/ul

response 223041

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	7.30	6.70
80.00	6.70	7.94
0.00	0.00	0.00

Quantitation Report (Qedit)

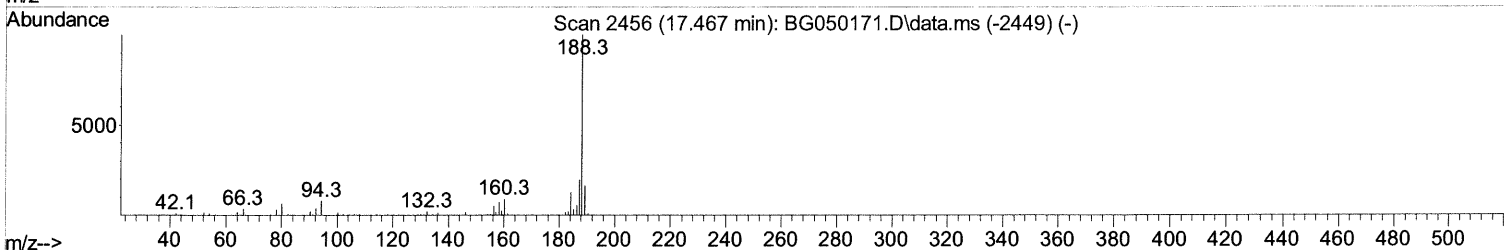
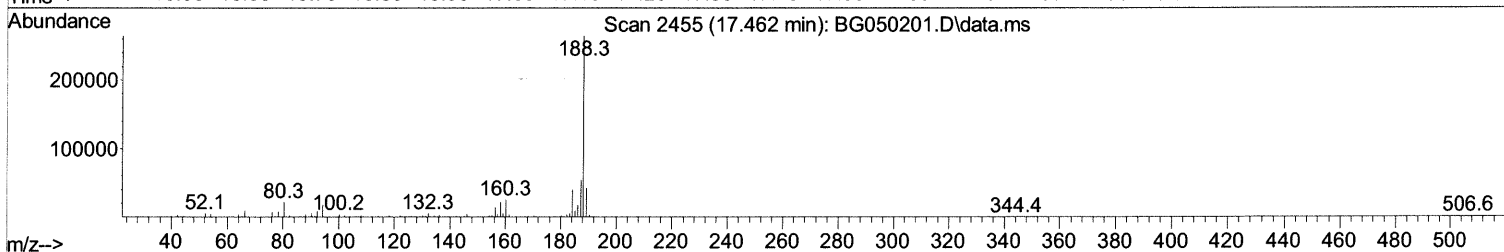
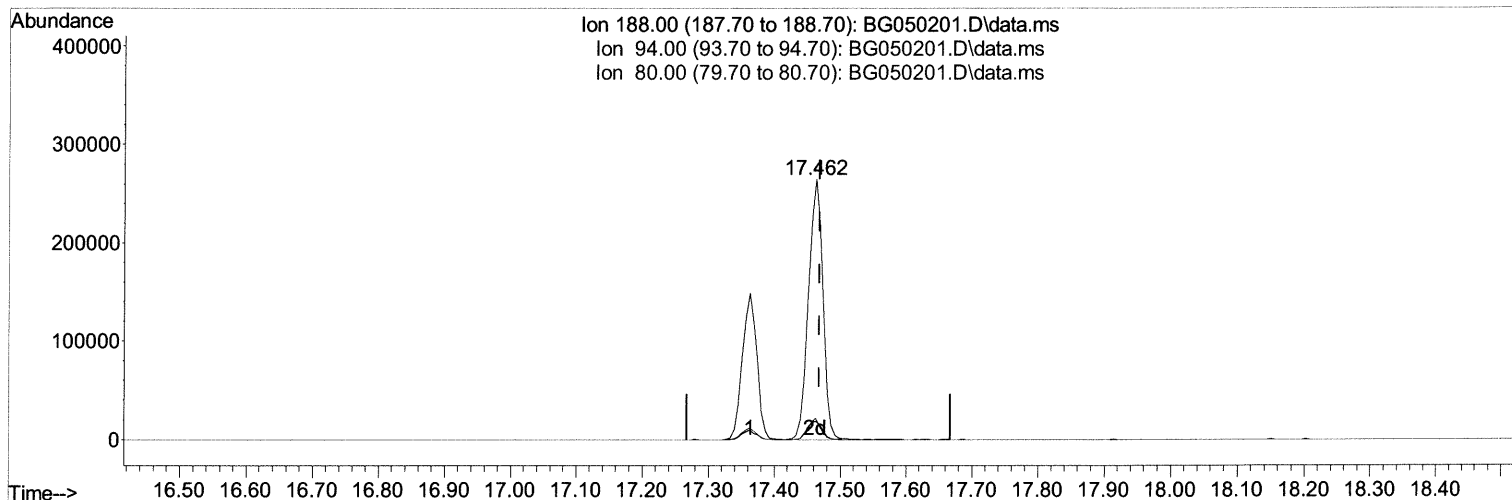
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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



TIC: BG050201.D\data.ms

(73) Anthracene-d10 (S)

17.462min (-0.005) 39.97 ng/ul m 09/10/21 JU

response 400255

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	7.30	6.96
80.00	6.70	8.15#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
 Data File : BG050201.D
 Acq On : 8 Sep 2021 15:54
 Operator : CG/JU
 Sample : M3608-15
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBKL1

Manual Integrations
 APPROVED

mohammad
 9/9/2021 8:41:22 AM

Quant Time: Sep 08 16:50:43 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 07 18:32:38 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.946	152	39385	20.000 ng/ul	-0.01
20) Naphthalene-d8	10.766	136	156474	20.000 ng/ul	-0.01
38) Acenaphthene-d10	14.607	164	101281	20.000 ng/ul	-0.01
64) Phenanthrene-d10	17.363	188	223041	20.000 ng/ul	# 0.00
79) Chrysene-d12	21.633	240	229477	20.000 ng/ul	0.00
88) Perylene-d12	24.847	264	240931	20.000 ng/ul	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.305	96	2088	2.621 ng/ul	0.00
4) Pyridine-d5	3.752	84	21385	7.562 ng/ul	0.01
7) Phenol-d5	7.129	99	27731	9.068 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.276	67	56312	31.531 ng/ul	0.00
11) 2-Chlorophenol-d4	7.482	132	64286	25.330 ng/ul	0.00
15) 4-Methylphenol-d8	8.674	113	46503	19.341 ng/ul	0.00
21) Nitrobenzene-d5	9.127	128	43229	40.250 ng/ul	0.00
24) 2-Nitrophenol-d4	9.849	143	43505	33.312 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.401	165	76946	30.870 ng/ul	0.00
31) 4-Chloroaniline-d4	10.907	131	86926	28.306 ng/ul	-0.01
46) Dimethylphthalate-d6	14.008	166	265547	34.723 ng/ul	-0.01
49) Acenaphthylene-d8	14.302	160	305364	33.534 ng/ul	-0.01
54) 4-Nitrophenol-d4	14.860	143	6747	7.312 ng/ul	0.00
60) Fluorene-d10	15.606	176	221980	34.958 ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.741	200	44389	32.993 ng/ul	-0.01
73) Anthracene-d10	17.462	188	400255m >	39.975 ng/ul >	0.00
81) Pyrene-d10	19.753	212	466483	40.811 ng/ul	0.00
92) Benzo(a)pyrene-d12	24.629	264	471379	36.959 ng/ul	-0.01

09/10/21 JU

Target Compounds Qvalue

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