

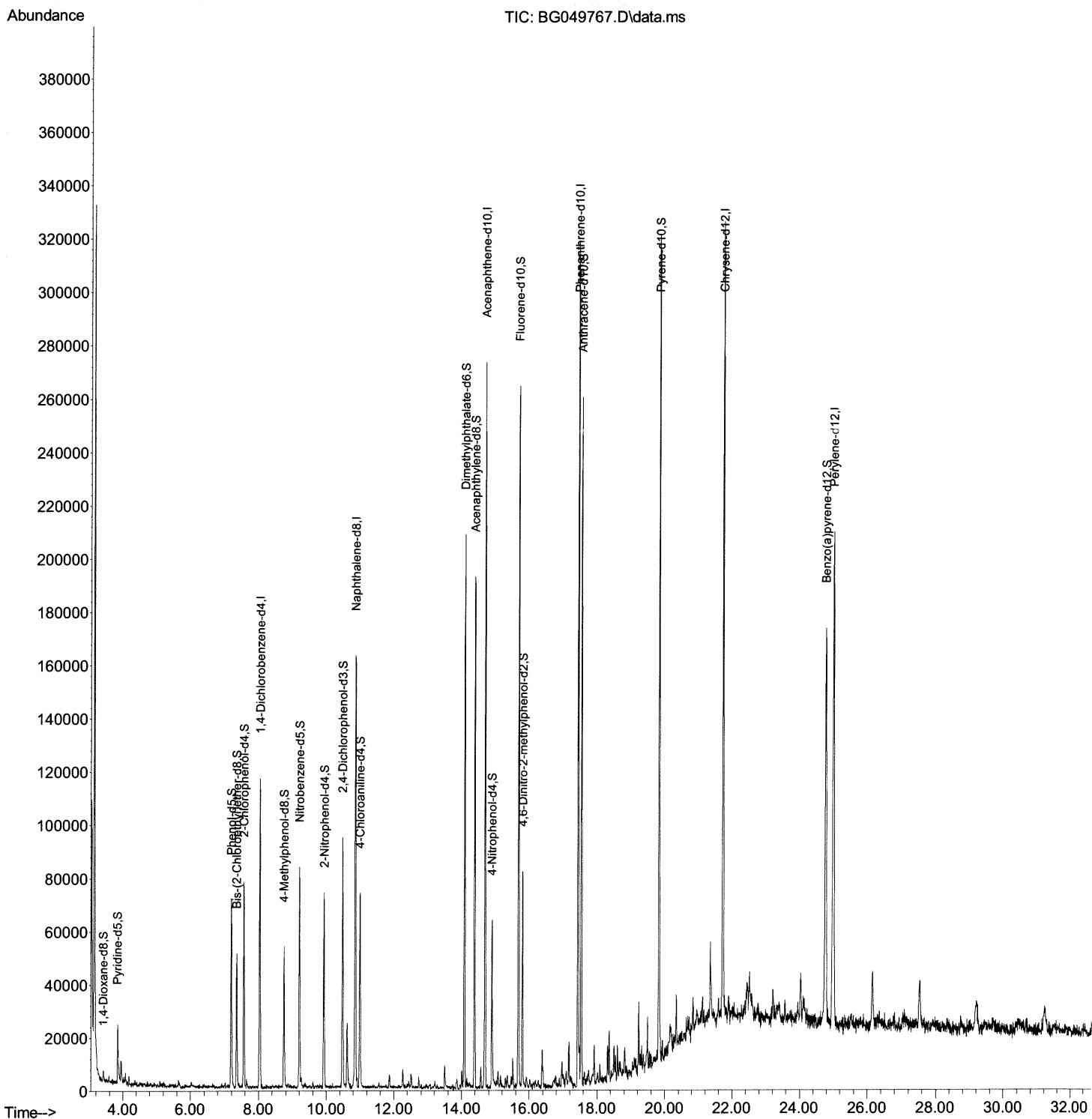
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081021\
Data File : BG049767.D
Acq On : 10 Aug 2021 21:56
Operator : CG/JU
Sample : M3182-15
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C00D3

Manual Integrations
APPROVED

mohammad
8/11/2021 9:33:08 PM

Quant Time: Aug 11 03:06:06 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)

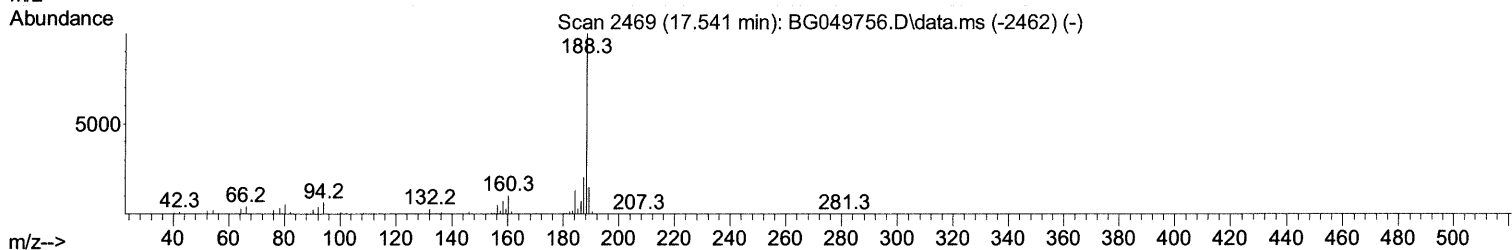
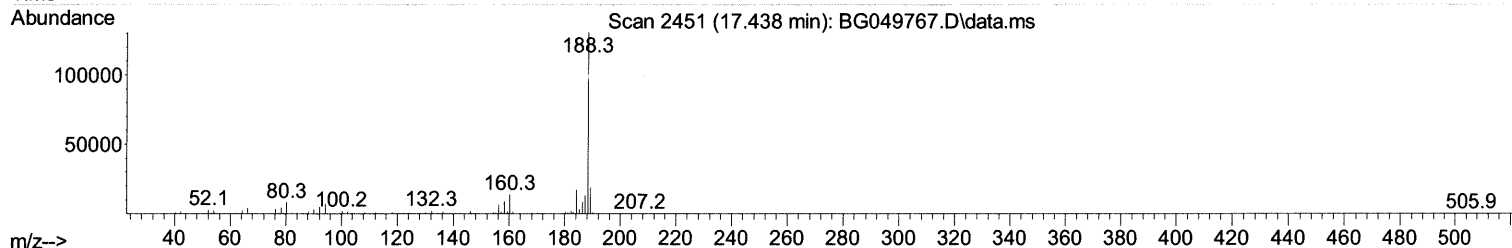
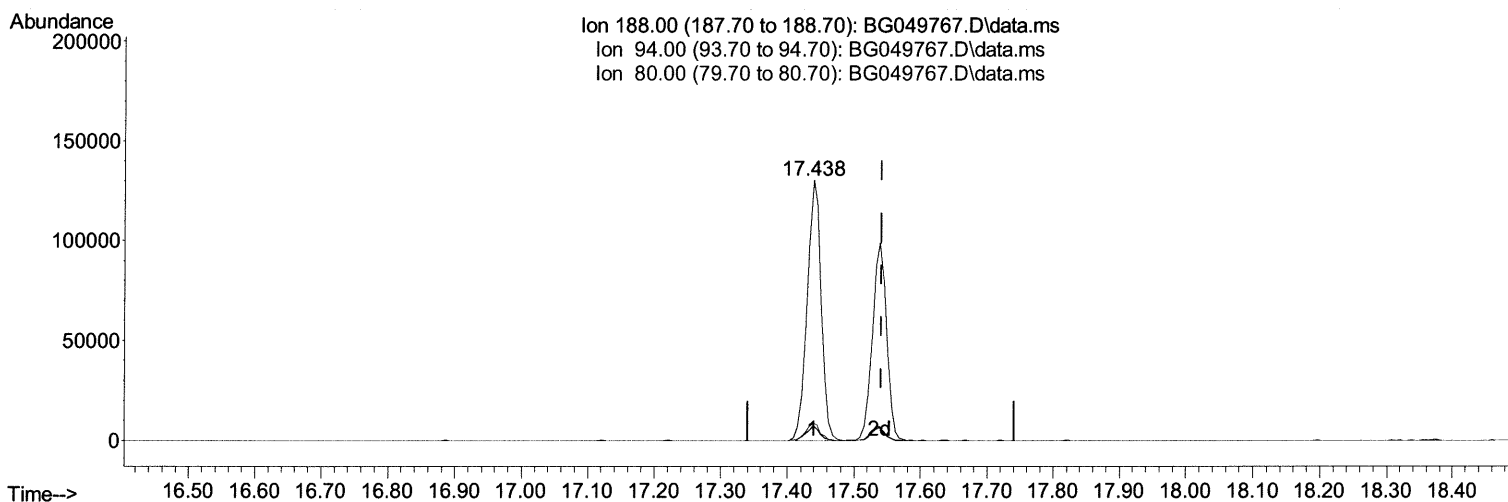
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TIC: BG049767.D\data.ms

(73) Anthracene-d10 (S)

17.438min (-0.102) 21.18 ng/ul

response 191728

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	5.60	5.65
80.00	6.00	6.42
0.00	0.00	0.00

Quantitation Report (Qedit)

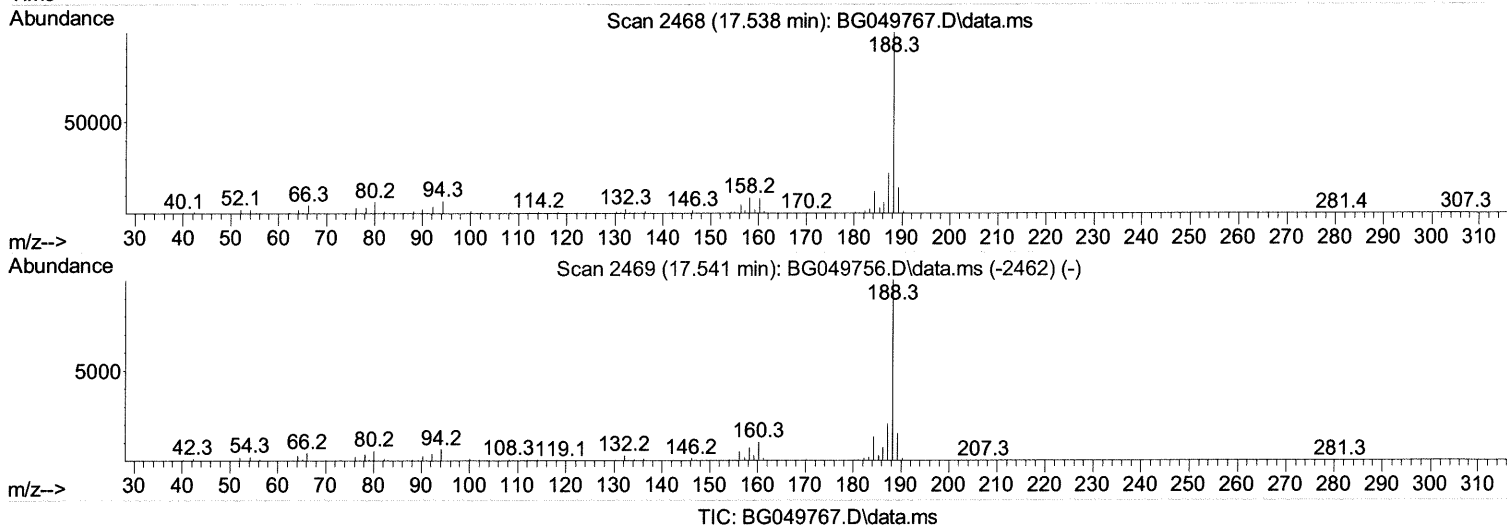
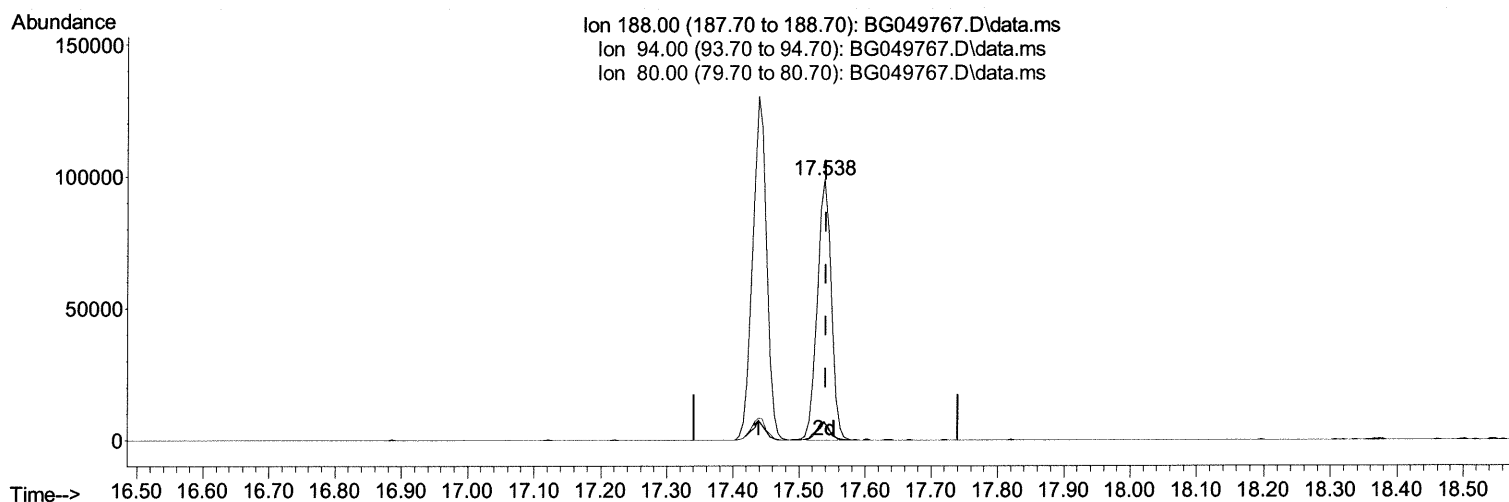
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Manual Integrations
APPROVED

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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



(73) Anthracene-d10 (S)

17.538min (-0.002) 16.17 ng/ul m 08/12/21 JU

response 146314

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	5.60	7.02#
80.00	6.00	6.47
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081021\
 Data File : BG049767.D
 Acq On : 10 Aug 2021 21:56
 Operator : CG/JU
 Sample : M3182-15
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleID :
 C00D3

Manual Integrations
 APPROVED

mohammad
 8/11/2021 9:33:08 PM

Quant Time: Aug 11 03:06:06 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

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.034	152	33636	20.000 ng/ul	0.00
20) Naphthalene-d8	10.853	136	135002	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.689	164	95325	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.438	188	191728	20.000 ng/ul	0.00
79) Chrysene-d12	21.721	240	174544	20.000 ng/ul	0.00
88) Perylene-d12	24.993	264	181240	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.416	96	1987	2.762 ng/ul	0.00
4) Pyridine-d5	3.845	84	13781	6.383 ng/ul	0.00
7) Phenol-d5	7.199	99	40084	13.131 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.352	67	25797	14.712 ng/ul	0.00
11) 2-Chlorophenol-d4	7.564	132	33489	15.328 ng/ul	0.00
15) 4-Methylphenol-d8	8.744	113	20702	8.894 ng/ul	0.00
21) Nitrobenzene-d5	9.203	128	17743	16.685 ng/ul	0.00
24) 2-Nitrophenol-d4	9.937	143	20751	17.086 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.477	165	33094	14.452 ng/ul	0.00
31) 4-Chloroaniline-d4	10.994	131	38779	12.370 ng/ul	0.00
46) Dimethylphthalate-d6	14.084	166	124506	16.644 ng/ul	0.00
49) Acenaphthylene-d8	14.378	160	144404	16.225 ng/ul	0.00
54) 4-Nitrophenol-d4	14.895	143	15169	12.542 ng/ul	0.00
60) Fluorene-d10	15.682	176	104725	16.286 ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.799	200	17468	13.537 ng/ul	0.00
73) Anthracene-d10	17.538	188	146314m	16.167 ng/ul	0.00
81) Pyrene-d10	19.823	212	175115	18.209 ng/ul	0.00
92) Benzo(a)pyrene-d12	24.764	264	160027	16.114 ng/ul	0.00

Target Compounds Qvalue

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