

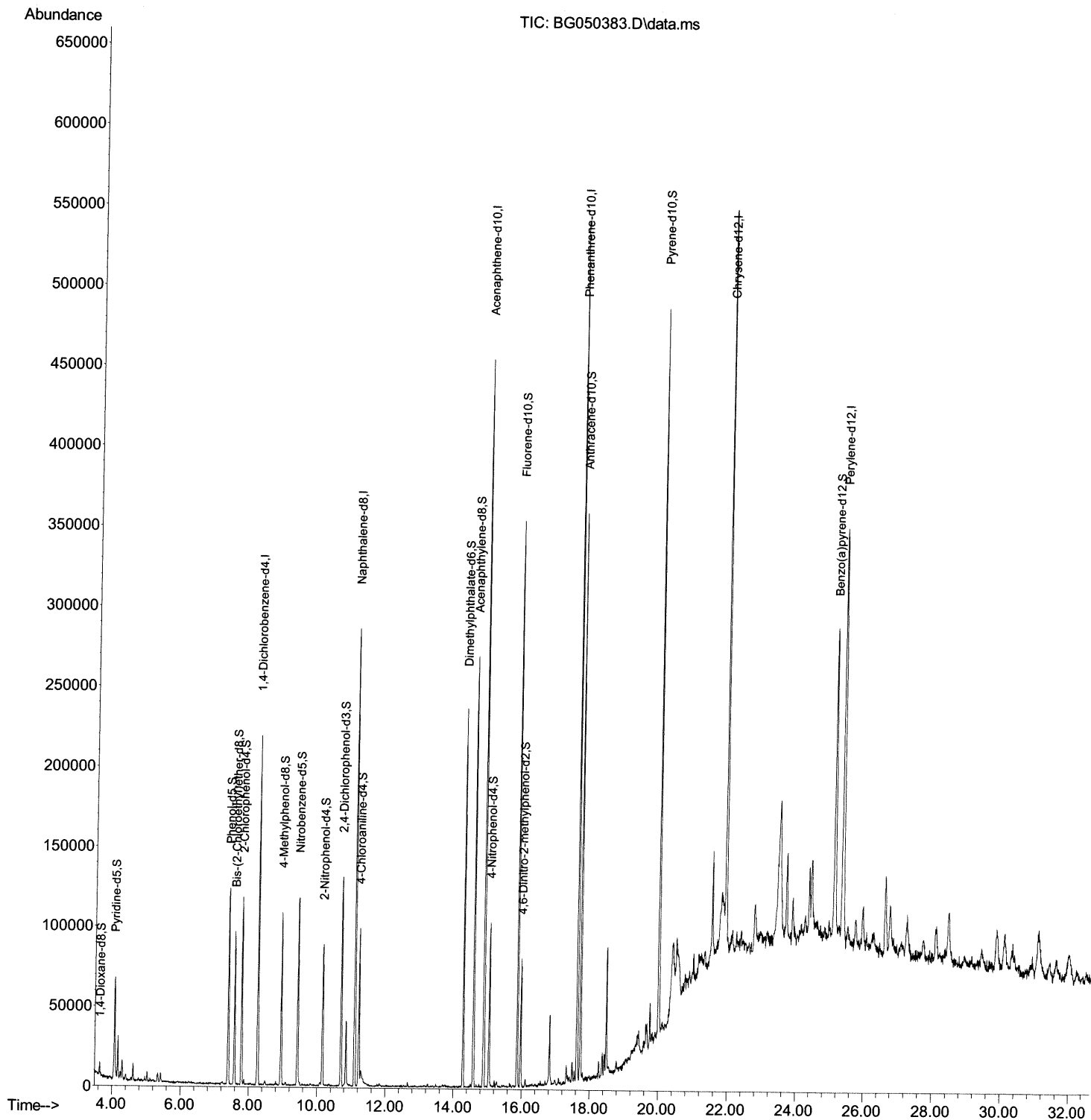
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
Data File : BG050383.D
Acq On : 2 Oct 2021 10:06
Operator : CG/JU
Sample : M3877-07
Misc :
ALS Vial : 68 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW2M3

Manual Integrations
APPROVED

mohammad
10/4/2021 11:51:03 AM

Quant Time: Oct 04 01:05:29 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.M
Quant Title : SVOA CALIBRATION
Qlast Update : Thu Sep 30 16:27:24 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

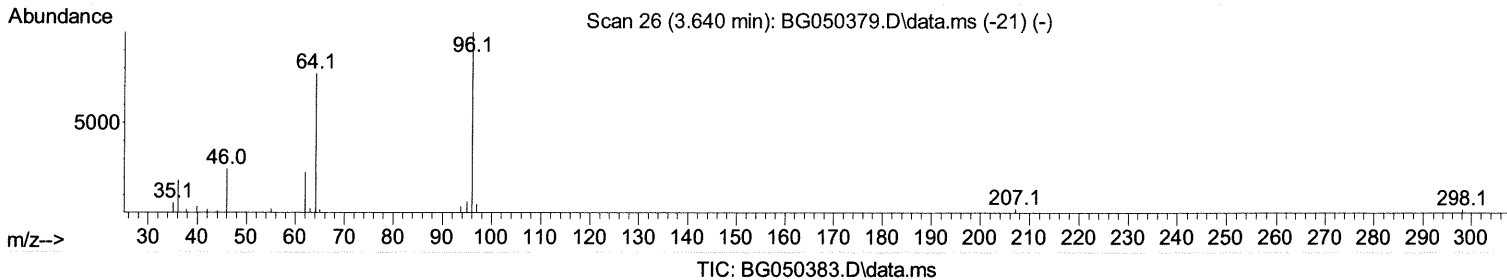
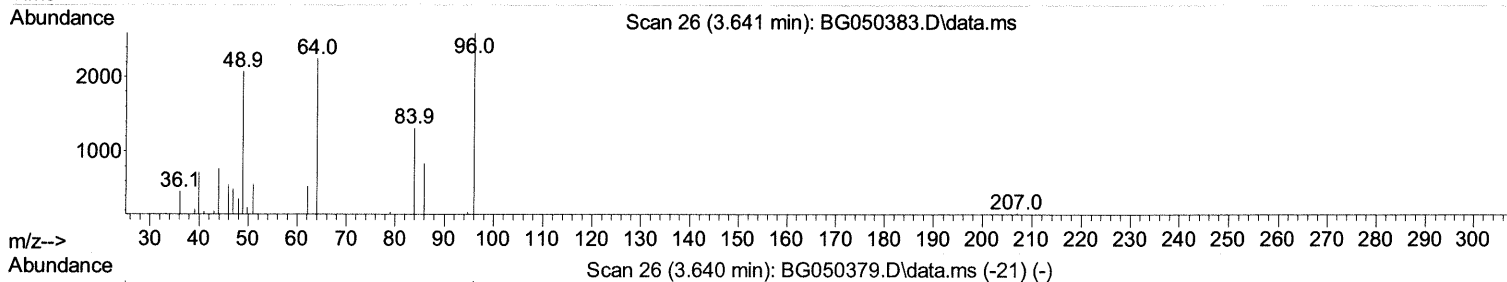
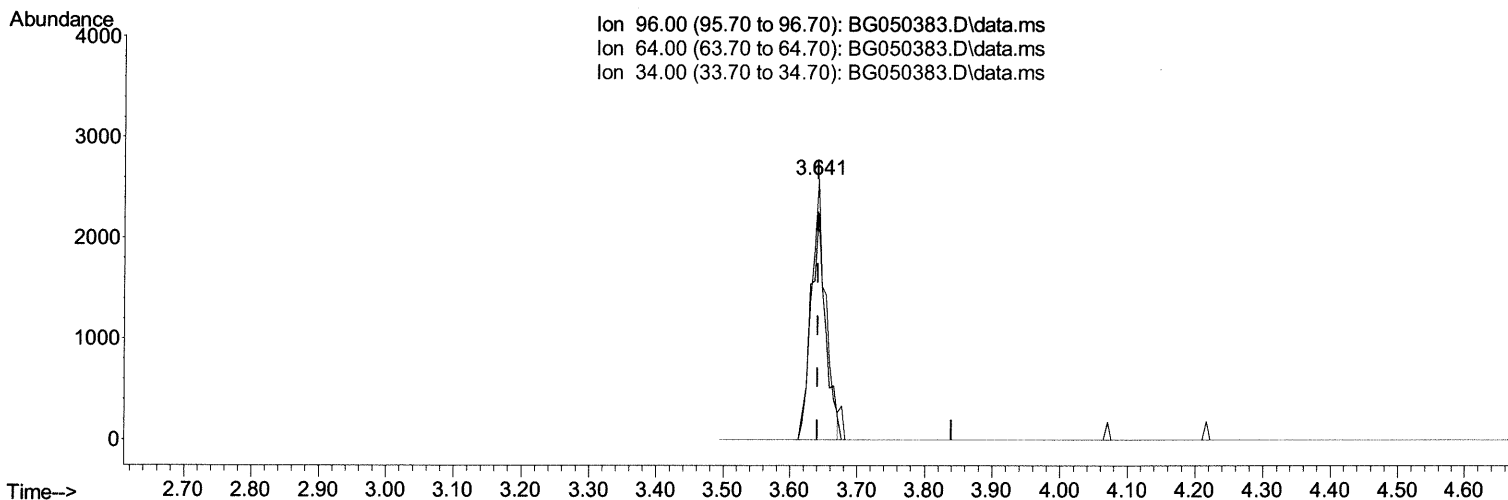
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(3) 1,4-Dioxane-d8 (S)

3.641min (+ 0.001) 2.49 ng/uL

response 3825

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	89.00	86.89
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

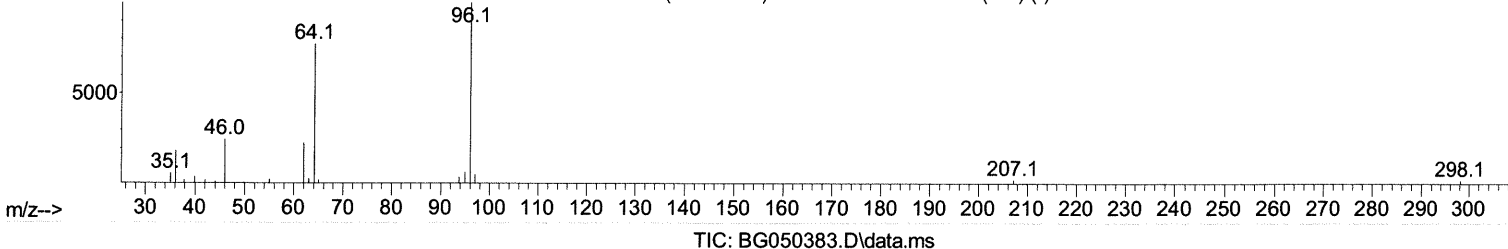
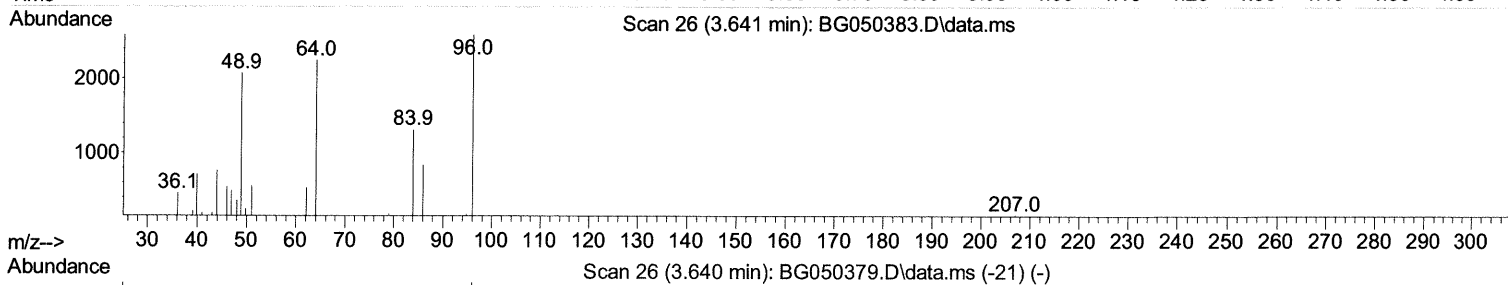
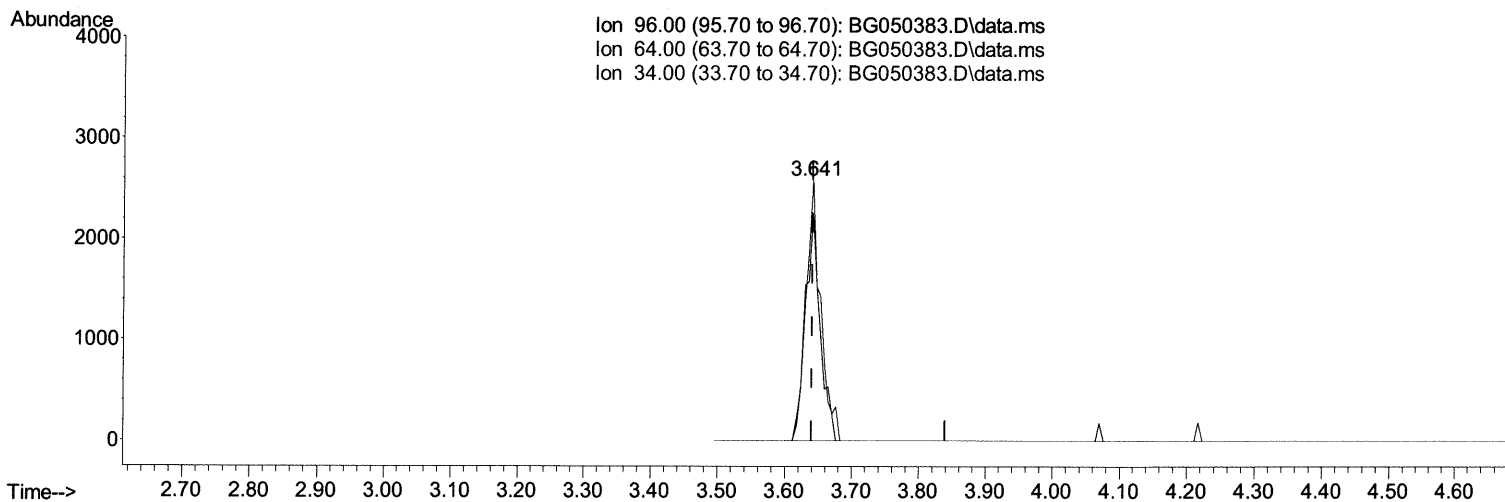
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TIC: BG050383.D\data.ms

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

3.641min (+ 0.001) 2.57 ng/uL m 10/05/21 JU

response 3942

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	89.00	86.89
34.00	0.00	0.00
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	8.271	152	56401	20.000	ng/ul	0.00
20) Naphthalene-d8	11.097	136	237164	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.886	164	153451	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.630	188	321579	20.000	ng/ul	0.00
79) Chrysene-d12	21.931	240	290390	20.000	ng/ul	-0.02
88) Perylene-d12	25.351	264	284056	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.641	96	3942m	2.568	ng/ul	> 0.00
4) Pyridine-d5	4.064	84	40688	8.895	ng/ul	0.00
7) Phenol-d5	7.401	99	70093	14.224	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.583	67	46735	15.070	ng/ul	0.00
11) 2-Chlorophenol-d4	7.795	132	49789	14.420	ng/ul	0.00
15) 4-Methylphenol-d8	8.958	113	42202	11.315	ng/ul	0.00
21) Nitrobenzene-d5	9.440	128	26165	15.770	ng/ul	0.00
24) 2-Nitrophenol-d4	10.163	143	27920	15.680	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.697	165	47856	13.610	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.220	131	51849	10.322	ng/ul	0.00
46) Dimethylphthalate-d6	14.281	166	152604	14.531	ng/ul	-0.01
49) Acenaphthylene-d8	14.581	160	192460	14.593	ng/ul	-0.01
54) 4-Nitrophenol-d4	15.045	143	23286	11.512	ng/ul	-0.01
60) Fluorene-d10	15.874	176	139349	14.728	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.973	200	18118	11.296	ng/ul	-0.01
73) Anthracene-d10	17.724	188	211399	15.561	ng/ul	-0.01
81) Pyrene-d10	19.998	212	246970	15.994	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.122	264	207015	14.494	ng/ul	-0.02

Target Compounds Qvalue

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