

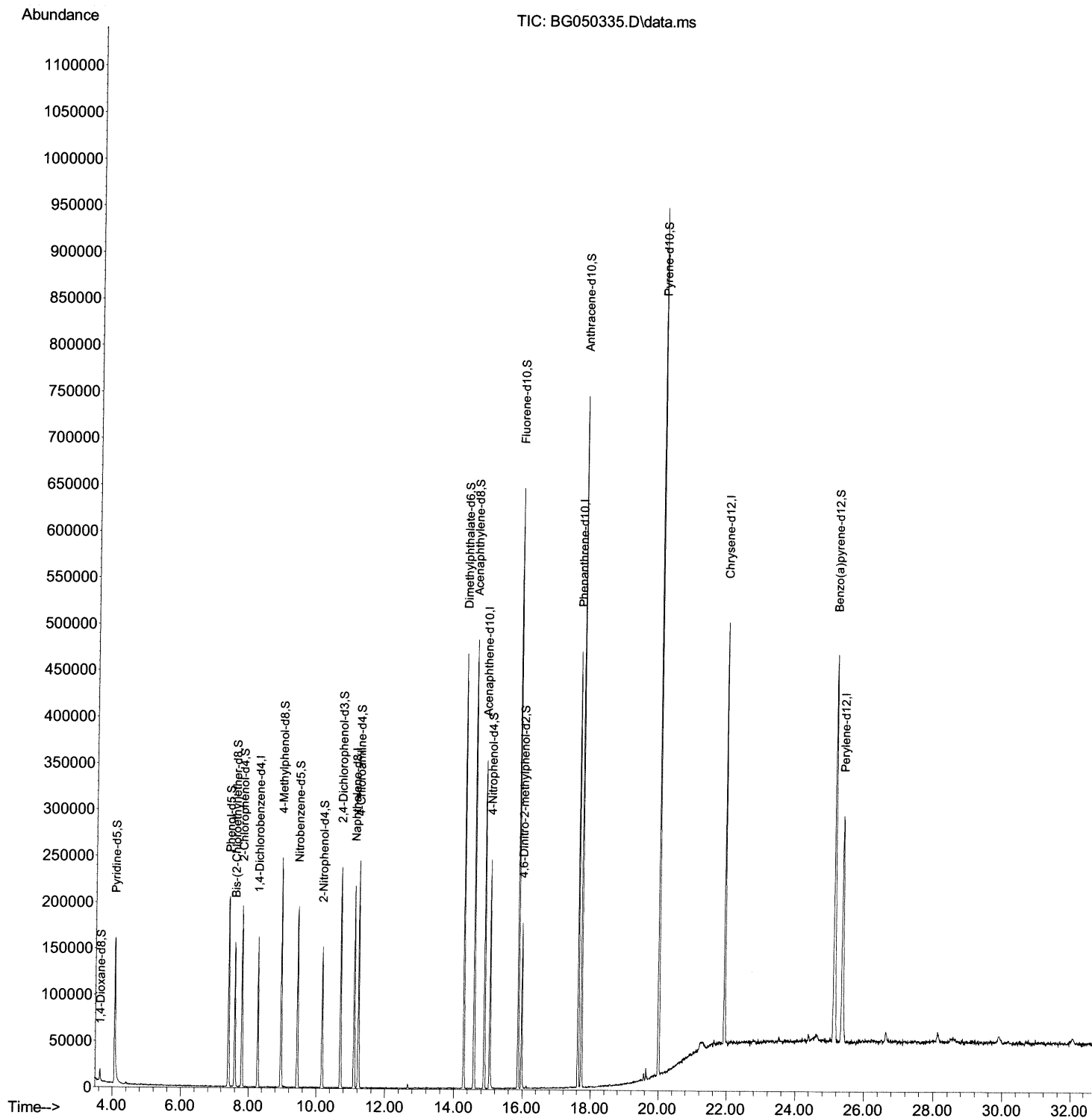
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
Data File : BG050335.D
Acq On : 30 Sep 2021 23:21
Operator : CG/JU
Sample : PB139240BL
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK240

Manual Integrations
APPROVED

mohammad
10/1/2021 11:30:16 AM

Quant Time: Oct 01 02:47:37 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)

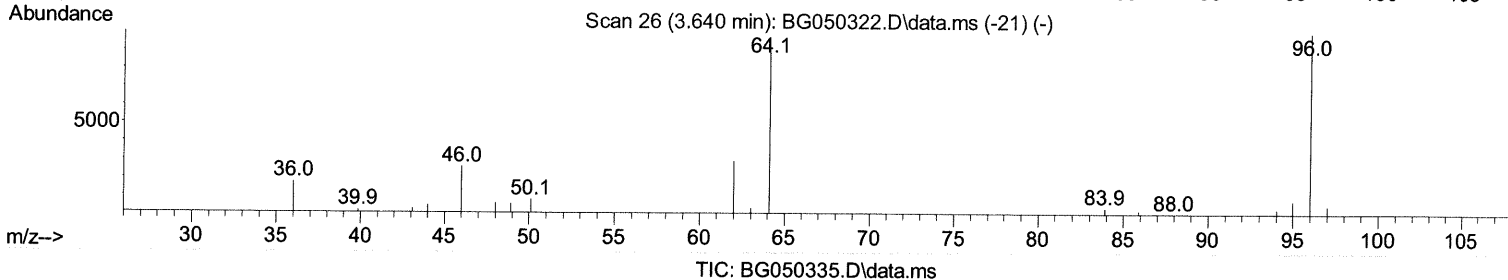
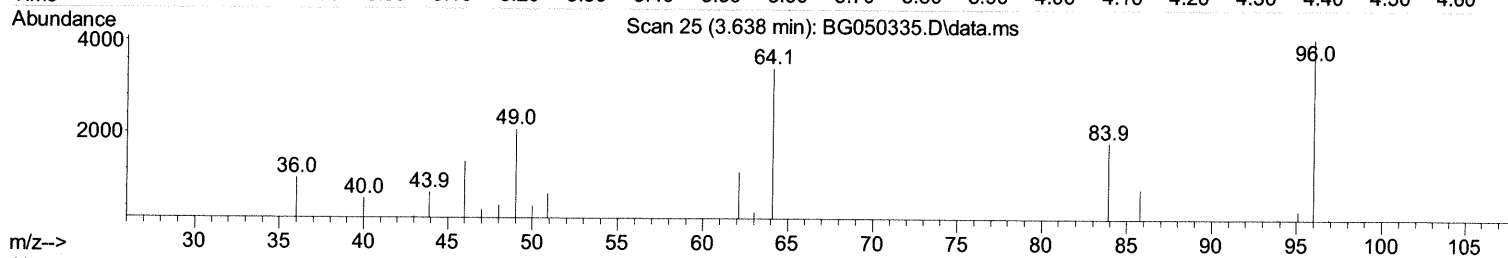
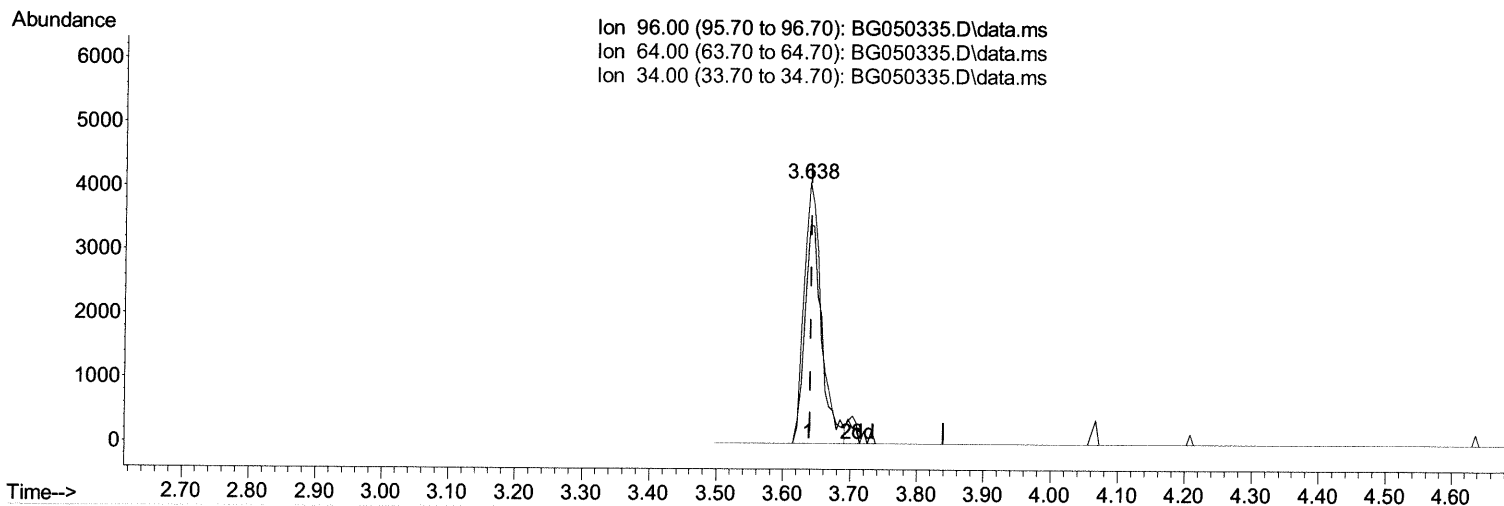
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(3) 1,4-Dioxane-d8 (S)

3.638min (-0.002) 6.00 ng/uL

response 7374

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	89.00	83.81
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

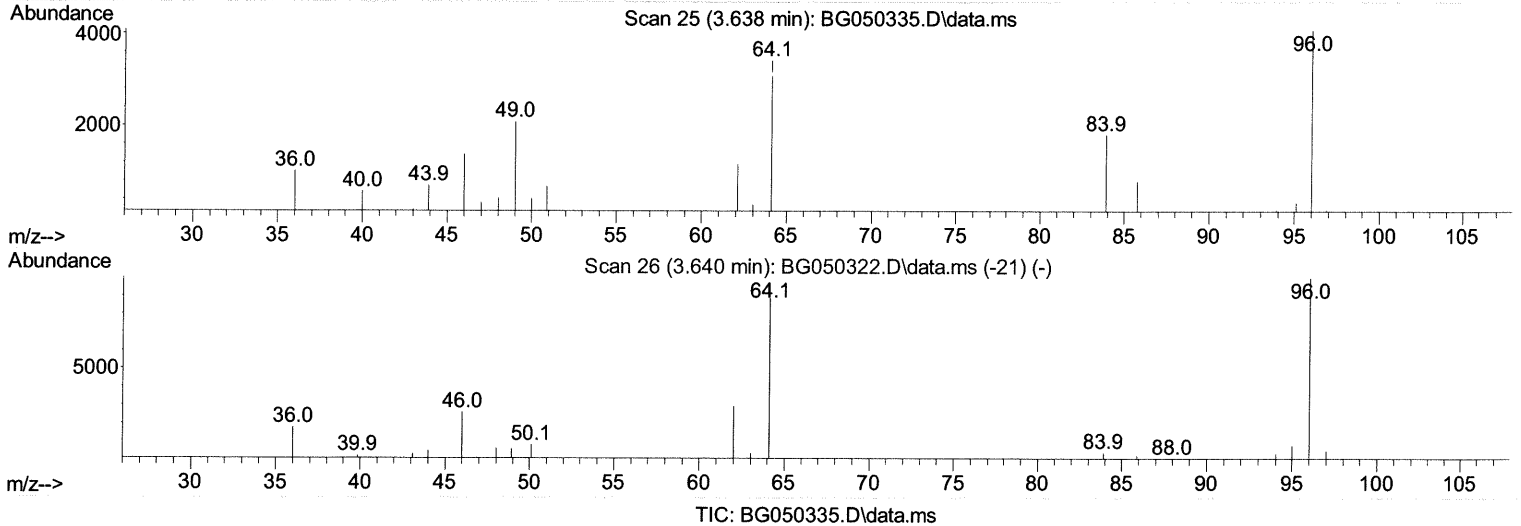
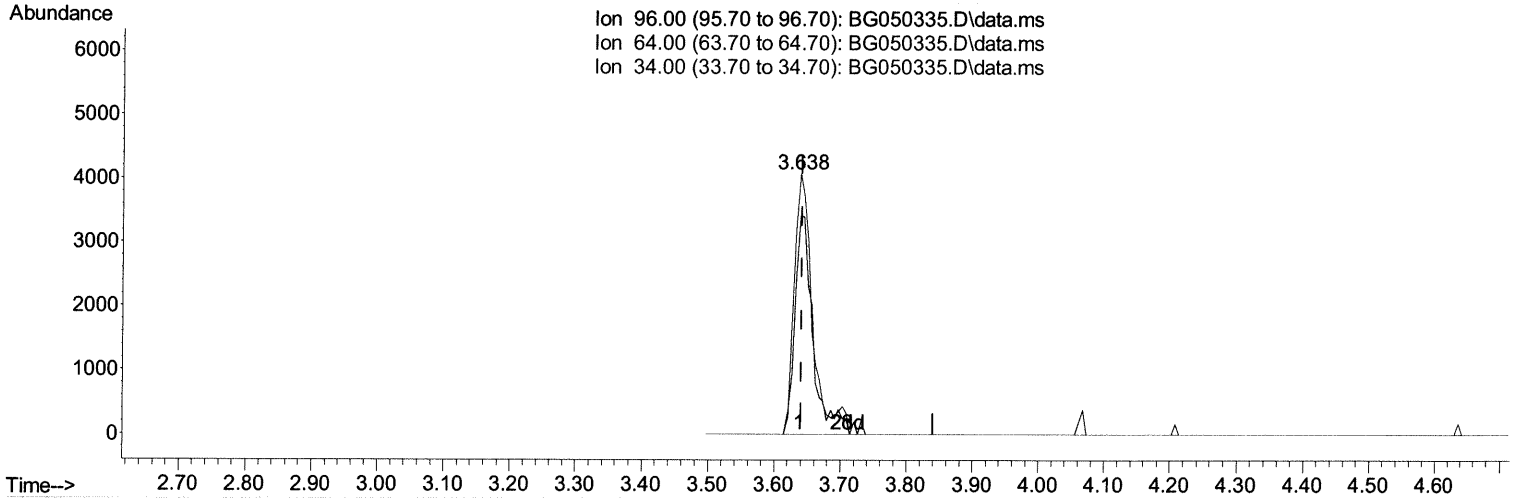
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(3) 1,4-Dioxane-d8 (S)

3.638min (-0.002) 6.32 ng/uL m 10/05/21 JU

response 7762

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	89.00	83.81
34.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.274	152	45125	20.000	ng/ul	0.00
20) Naphthalene-d8	11.100	136	183232	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.889	164	121838	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.633	188	280953	20.000	ng/ul	0.00
79) Chrysene-d12	21.934	240	282373	20.000	ng/ul	-0.01
88) Perylene-d12	25.360	264	264735	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.638	96	7762m>	6.320	ng/ul>	0.00 10/05/21 JU
4) Pyridine-d5	4.067	84	108644	29.685	ng/ul	0.00
7) Phenol-d5	7.404	99	124634	31.613	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.586	67	82263	33.154	ng/ul	0.00
11) 2-Chlorophenol-d4	7.798	132	88243	31.944	ng/ul	0.00
15) 4-Methylphenol-d8	8.961	113	95270	31.926	ng/ul	0.00
21) Nitrobenzene-d5	9.443	128	45616	35.587	ng/ul	0.00
24) 2-Nitrophenol-d4	10.166	143	46774	34.000	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.700	165	83910	30.888	ng/ul	0.00
31) 4-Chloroaniline-d4	11.223	131	125332	32.295	ng/ul	0.00
46) Dimethylphthalate-d6	14.284	166	296922	35.609	ng/ul	0.00
49) Acenaphthylene-d8	14.584	160	352796	33.691	ng/ul	0.00
54) 4-Nitrophenol-d4	15.048	143	53188	33.119	ng/ul	0.00
60) Fluorene-d10	15.877	176	256401	34.130	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.976	200	40761	29.088	ng/ul	0.00
73) Anthracene-d10	17.733	188	437177	36.833	ng/ul	0.00
81) Pyrene-d10	20.007	212	539568	35.936	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.125	264	452889	34.022	ng/ul	-0.01

Target Compounds Qvalue

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