Quantitation Report (QT Reviewed)

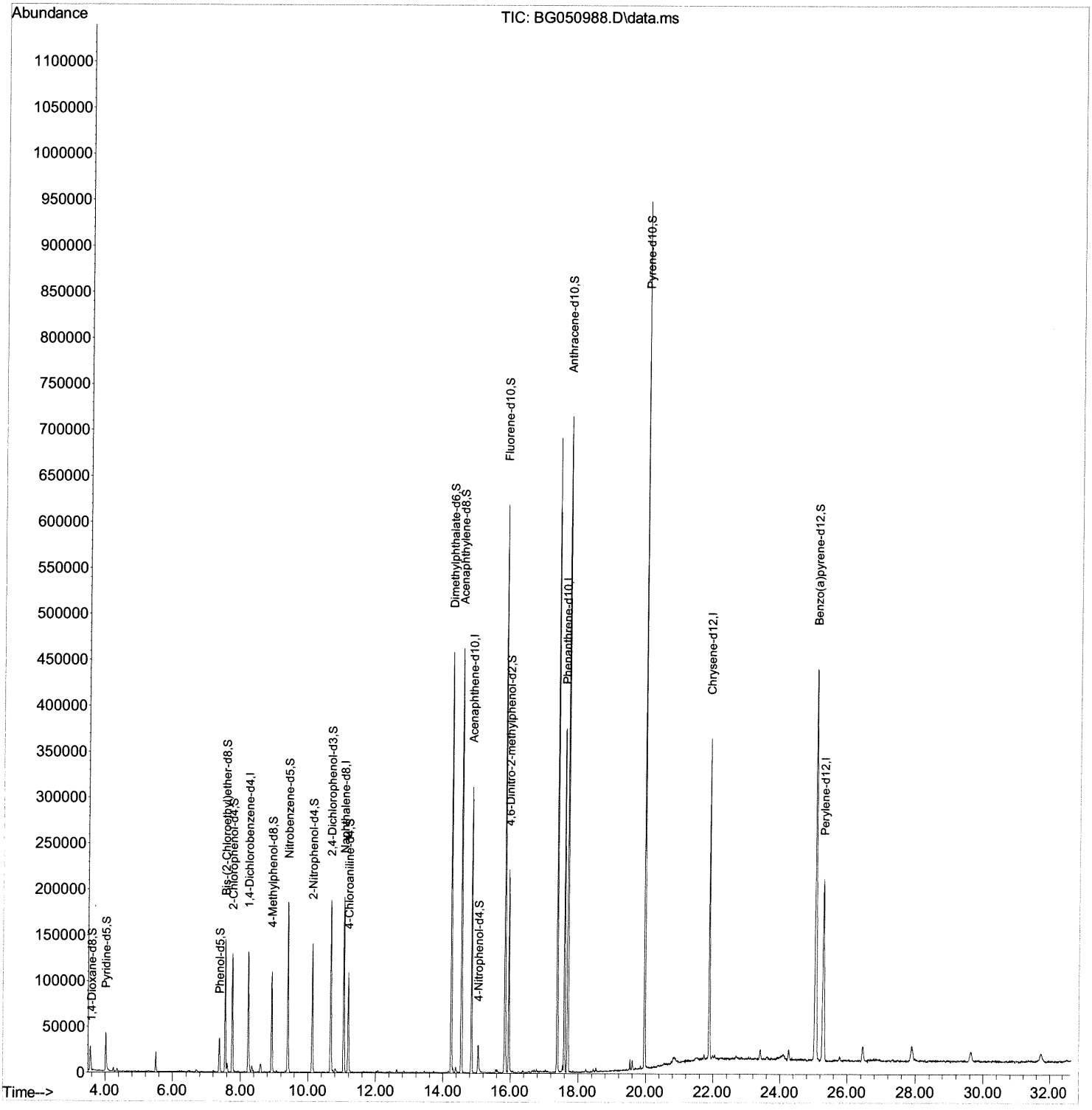
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050988.D
Acq On : 12 Nov 2021 5:23
Operator : CG/JU
Sample : M4542-09
Misc :
ALS Vial : 27 Sample Multiplier: 1

Quant Time: Nov 12 06:05:58 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Nov 11 12:40:48 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
BGGF3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/12/2021
Supervised By :mohammad ahmed 11/17/2021



Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050988.D
Acq On : 12 Nov 2021 5:23
Operator : CG/JU
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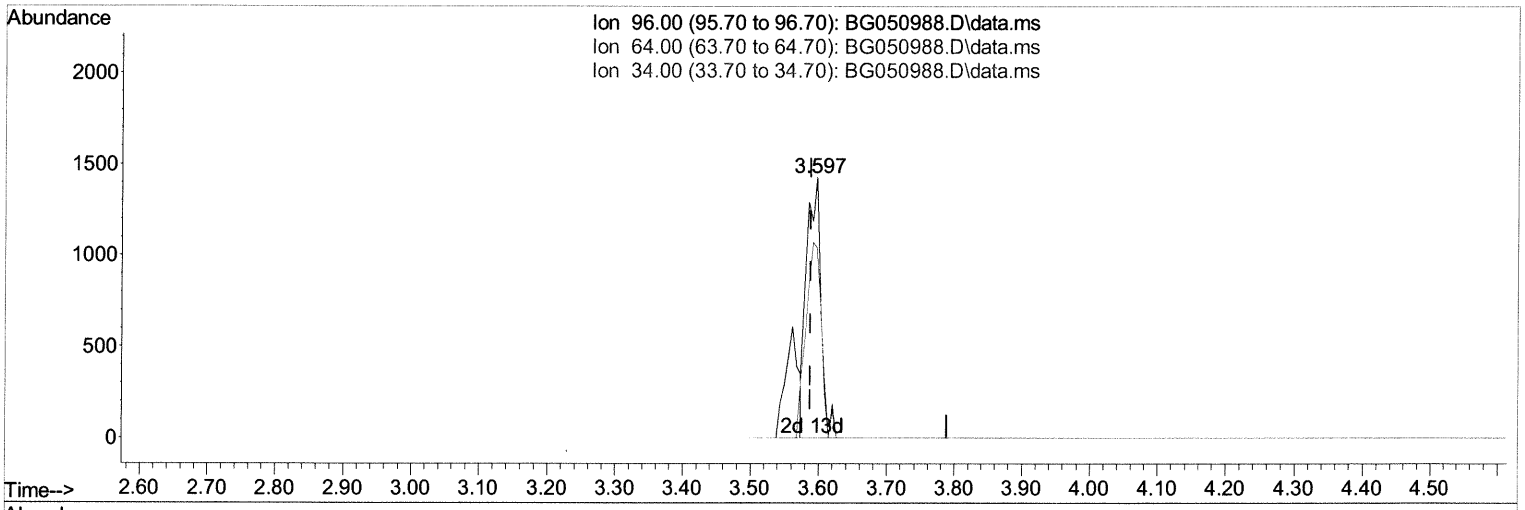

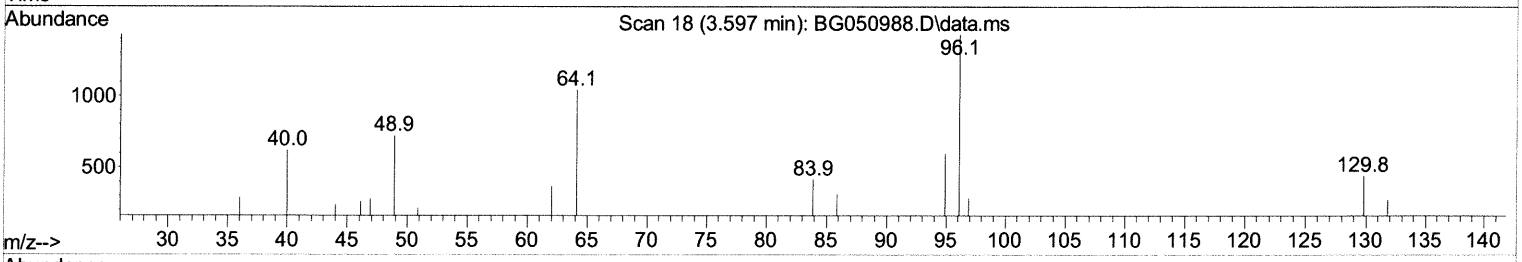

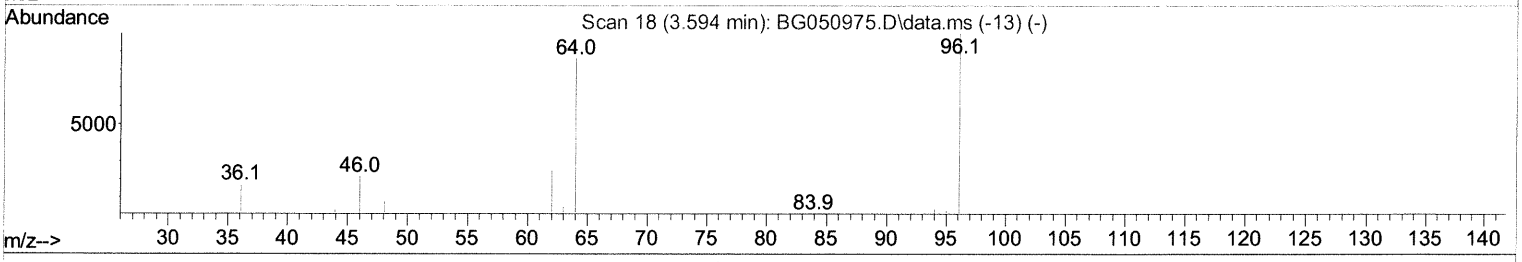

TIC: BG050988.D\data.ms

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

3.597min (+ 0.009) 1.90 ng/uL

response 2024

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	72.90
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/12/2021
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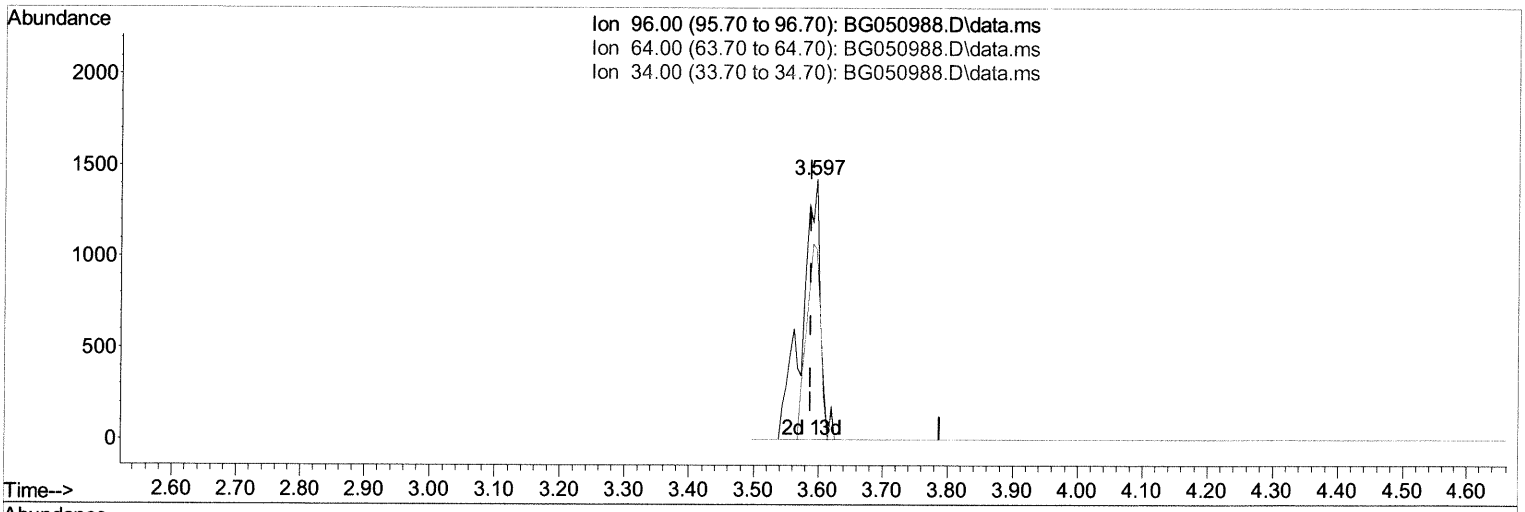

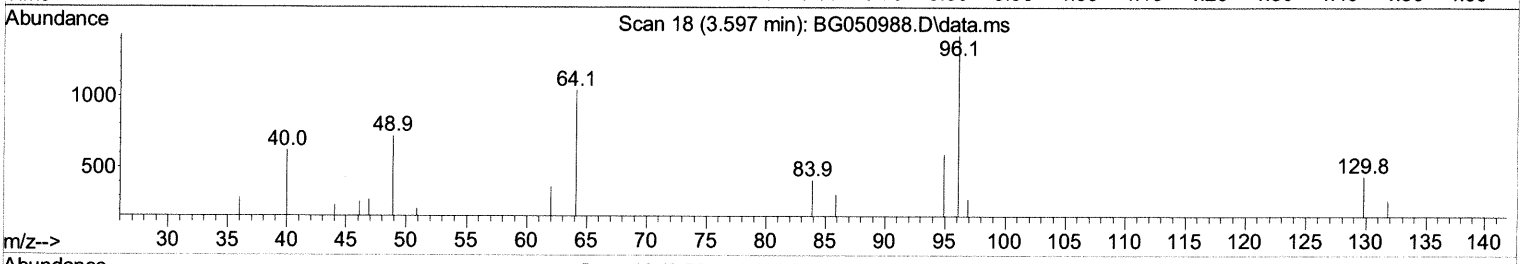

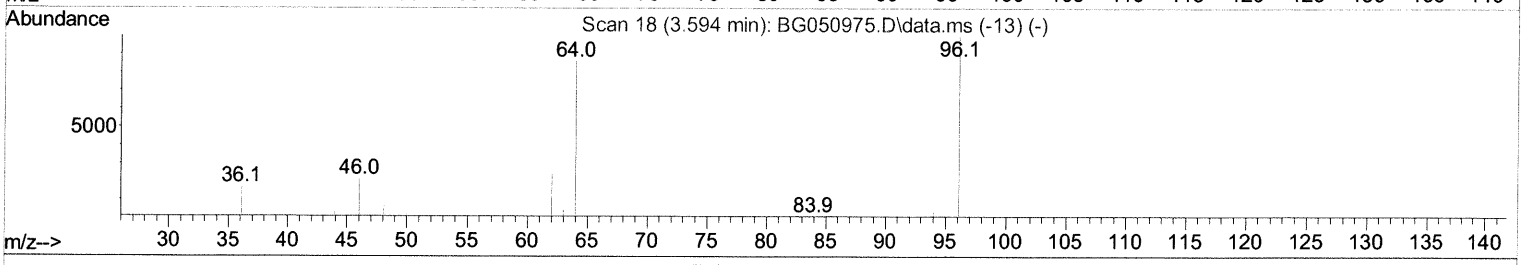

TIC: BG050988.D\data.ms

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

3.597min (+ 0.009) 2.70 ng/uL m 11/13/21 JU

response 2882

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	72.90
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
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Instrument :
BNA_G
ClientSampleId :
BGGF3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/12/2021
Supervised By :mohammad ahmed 11/17/2021

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.226	152	34409	20.00	ng/ul	0.00
20) Naphthalene-d8	11.053	136	157421	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.848	164	107709	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.598	188	235490	20.00	ng/ul	0.00
79) Chrysene-d12	21.893	240	212086	20.00	ng/ul	0.00
88) Perylene-d12	25.289	264	205778	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.597	96	2882m	> 2.70	ng/uL	> 0.00
4) Pyridine-d5	4.020	84	28298	8.87	ng/ul	0.00
7) Phenol-d5	7.375	99	22975	6.26	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.545	67	75060	31.65	ng/ul	0.00
11) 2-Chlorophenol-d4	7.756	132	59443	23.37	ng/ul	0.00
15) 4-Methylphenol-d8	8.926	113	43655	15.11	ng/ul	0.00
21) Nitrobenzene-d5	9.402	128	43669	32.64	ng/ul	0.00
24) 2-Nitrophenol-d4	10.124	143	46562	31.30	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.665	165	69811	27.86	ng/ul	0.00
31) 4-Chloroaniline-d4	11.188	131	61362	16.17	ng/ul	0.00
46) Dimethylphthalate-d6	14.243	166	292075	35.44	ng/ul	0.00
49) Acenaphthylene-d8	14.549	160	339954	33.11	ng/ul	0.00
54) 4-Nitrophenol-d4	15.048	143	8822	5.90	ng/ul	0.01
60) Fluorene-d10	15.841	176	244999	33.56	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.953	200	47843	33.51	ng/ul	0.00
73) Anthracene-d10	17.692	188	423881	38.07	ng/ul	0.00
81) Pyrene-d10	19.972	212	506616	36.98	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.054	264	430123	37.81	ng/ul	0.00

11/13/21 JU

Target Compounds Qvalue

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