

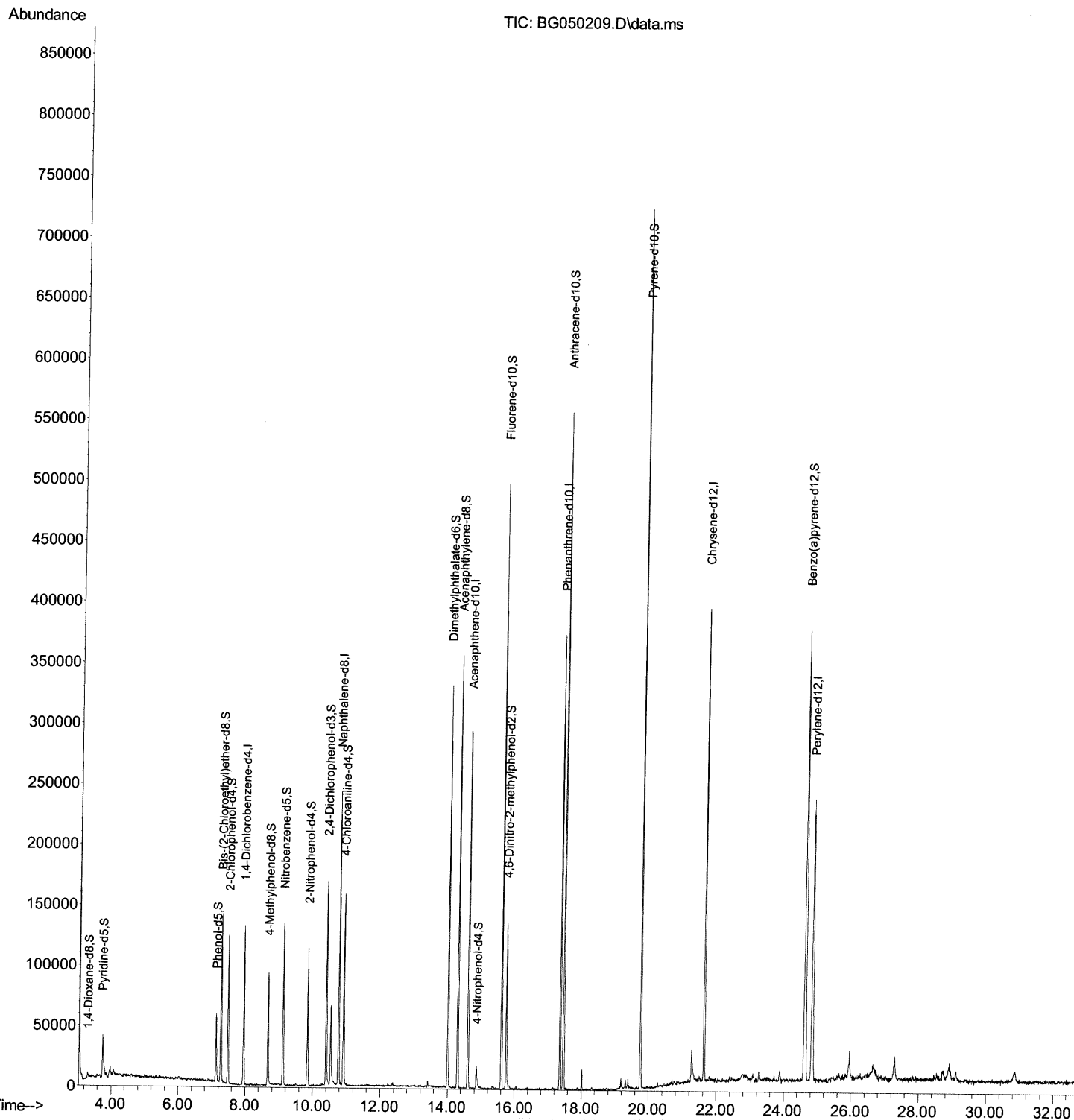
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050209.D
Acq On : 8 Sep 2021 22:08
Operator : CG/JU
Sample : M3608-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKK3

Manual Integrations
APPROVED

mohammad
9/9/2021 11:46:30 AM

Quant Time: Sep 09 00:57:15 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)

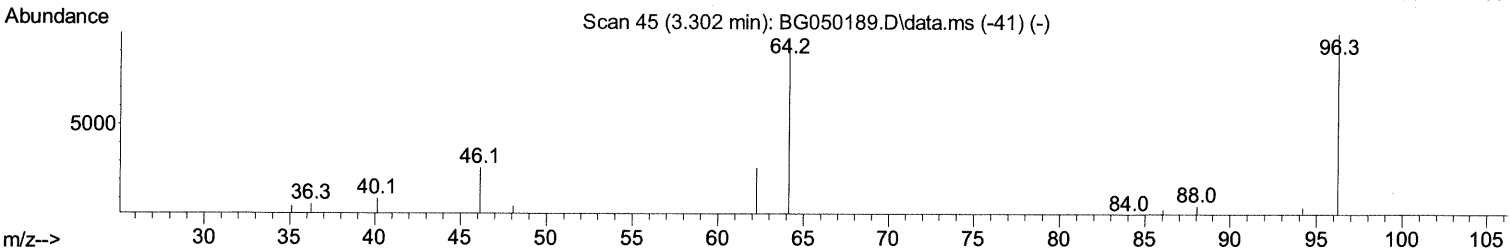
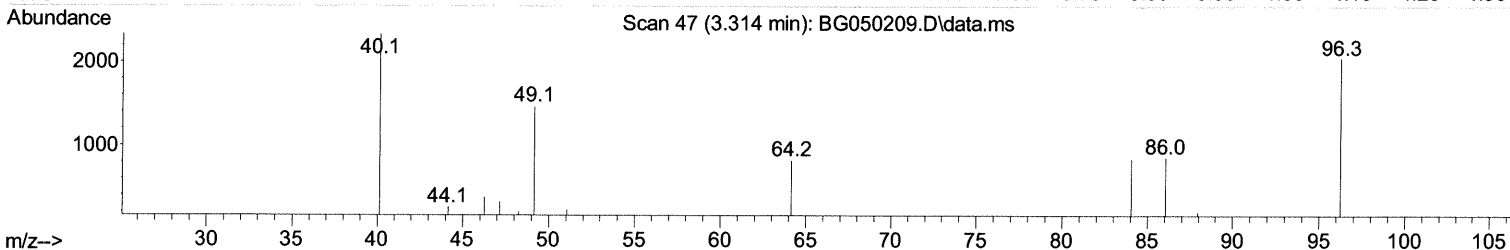
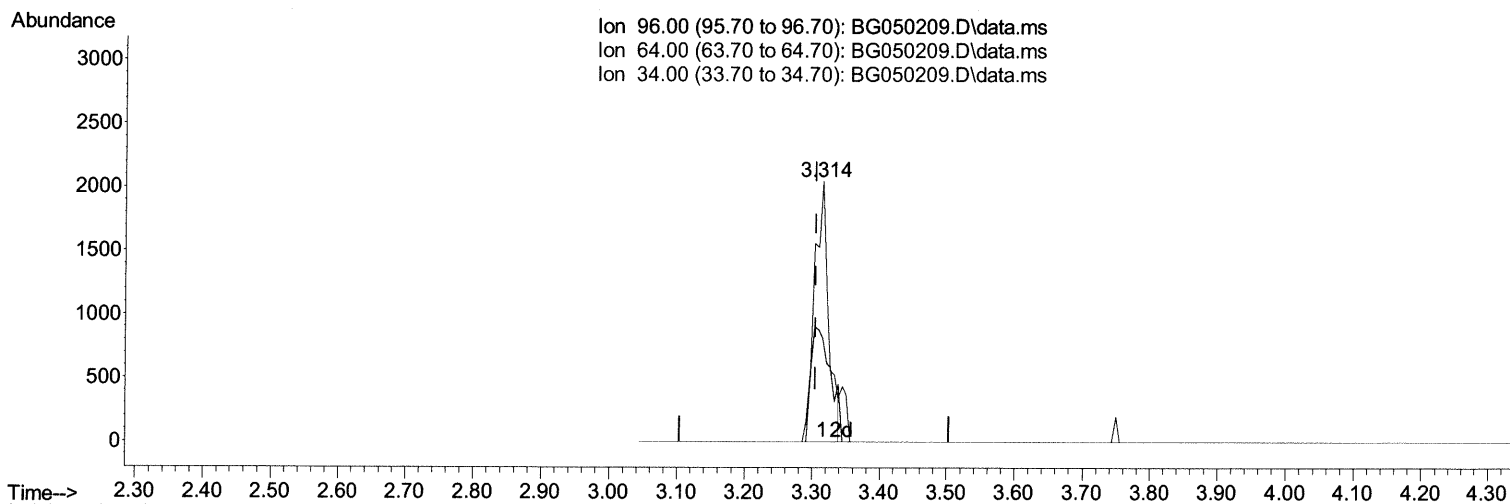
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TIC: BG050209.D\data.ms

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

3.314min (+ 0.010) 3.81 ng/uL

response 2984

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	63.40	39.68#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

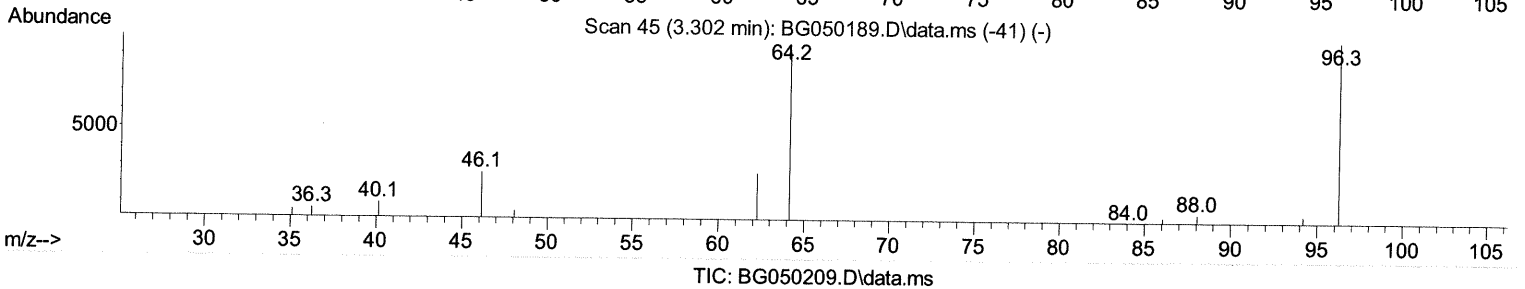
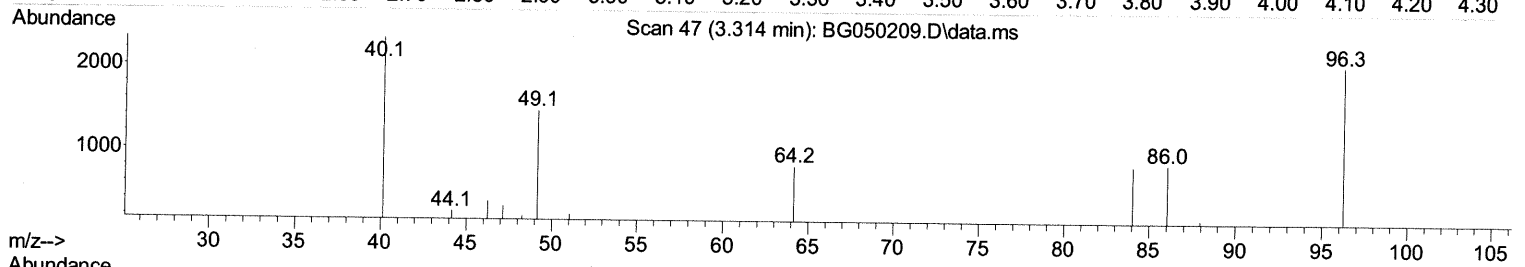
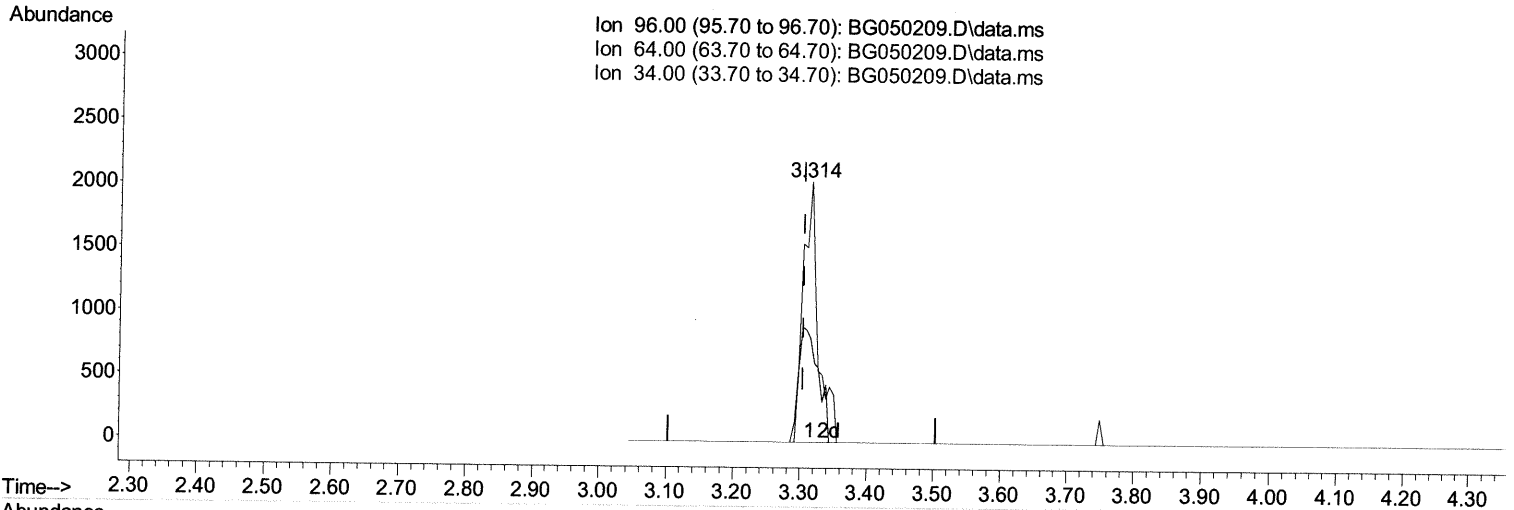
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TIC: BG050209.D\data.ms

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

3.314min (+ 0.010) 4.17 ng/uL m *09/10/21 JU*

response 3265

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	63.40	39.68#
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	7.949	152	38728	20.000	ng/u1	0.00
20) Naphthalene-d8	10.763	136	200235	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.611	164	108094	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.360	188	238048	20.000	ng/u1	0.00
79) Chrysene-d12	21.631	240	251565	20.000	ng/u1	0.00
88) Perylene-d12	24.844	264	247936	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.314	96	3265m>	4.168	ng/u1	> 0.01
4) Pyridine-d5	3.743	84	26624	9.575	ng/u1	0.00
7) Phenol-d5	7.127	99	37063	12.326	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.268	67	69347	39.489	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.479	132	59632	23.895	ng/u1	0.00
15) 4-Methylphenol-d8	8.678	113	37753	15.968	ng/u1	0.00
21) Nitrobenzene-d5	9.124	128	31380	22.832	ng/u1	0.00
24) 2-Nitrophenol-d4	9.847	143	36525	21.855	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	70014	21.950	ng/u1	0.00
31) 4-Chloroaniline-d4	10.904	131	89615	22.804	ng/u1	-0.01
46) Dimethylphthalate-d6	14.006	166	239946	29.398	ng/u1	-0.01
49) Acenaphthylene-d8	14.300	160	258538	26.602	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.864	143	5802	5.891	ng/u1	0.01
60) Fluorene-d10	15.604	176	195421	28.836	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	34259	23.858	ng/u1	0.00
73) Anthracene-d10	17.460	188	346122	32.389	ng/u1	0.00
81) Pyrene-d10	19.751	212	416054	33.203	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.627	264	403826	30.768	ng/u1	-0.01

Target Compounds Qvalue

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