

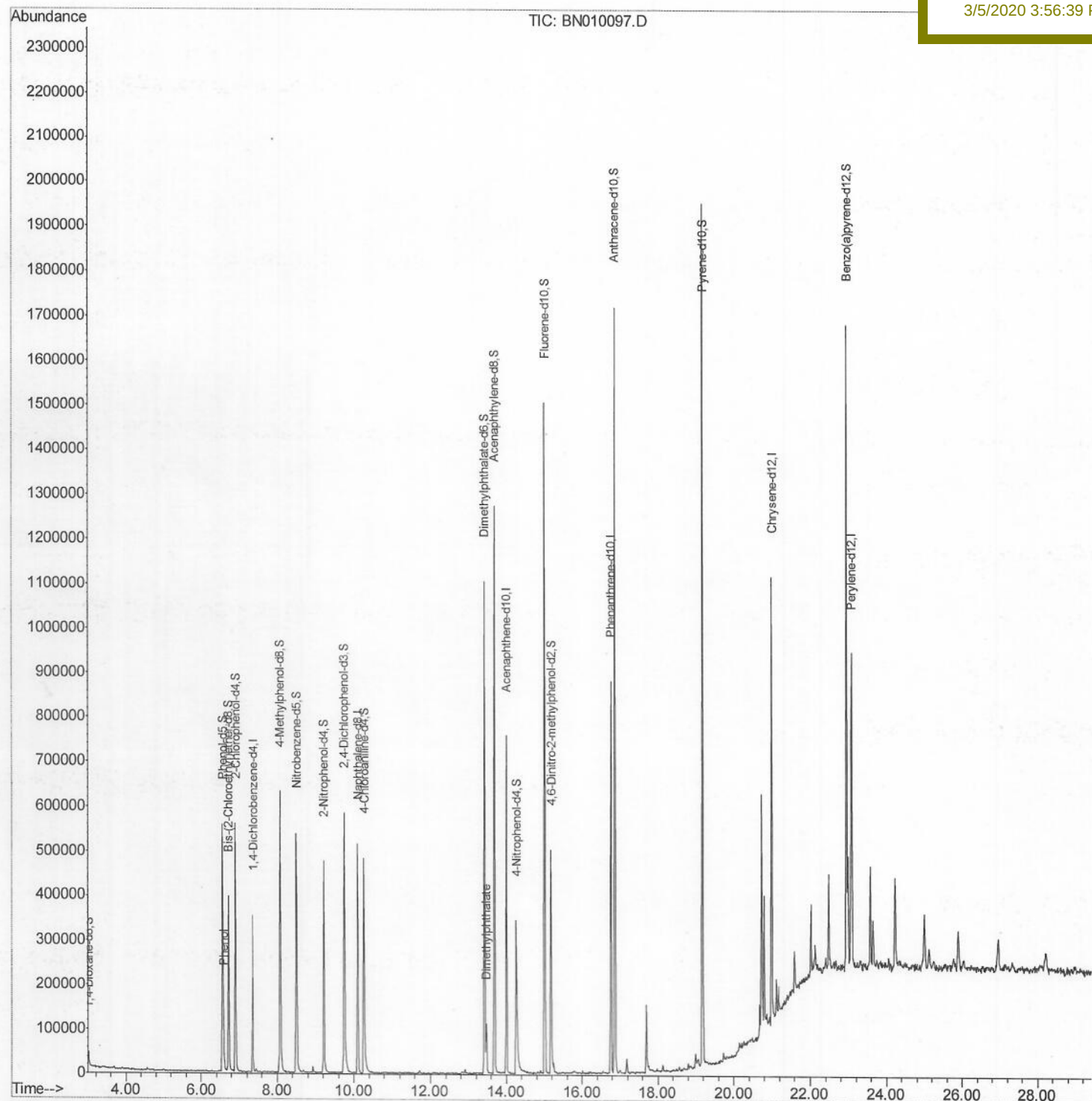
Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030420\
Data File : BN010097.D
Acq On : 05 Mar 2020 02:27
Operator : CG/JU
Sample : L1617-11
Misc :
ALS Vial : 24 Sample Multiplier: 1

Quant Time: Mar 05 08:16:43 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Mar 04 01:30:16 2020
Response via : Initial Calibration

Instrument :
BNA_N
ClientSampleId :
DBDS5

Manual Integrations
APPROVED

mohammad
3/5/2020 3:56:39 PM



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN030420\
Quantitation Report (Qedit)

Data File : BN010097.D

Acq On : 05 Mar 2020 02:27

Operator : CG/JU

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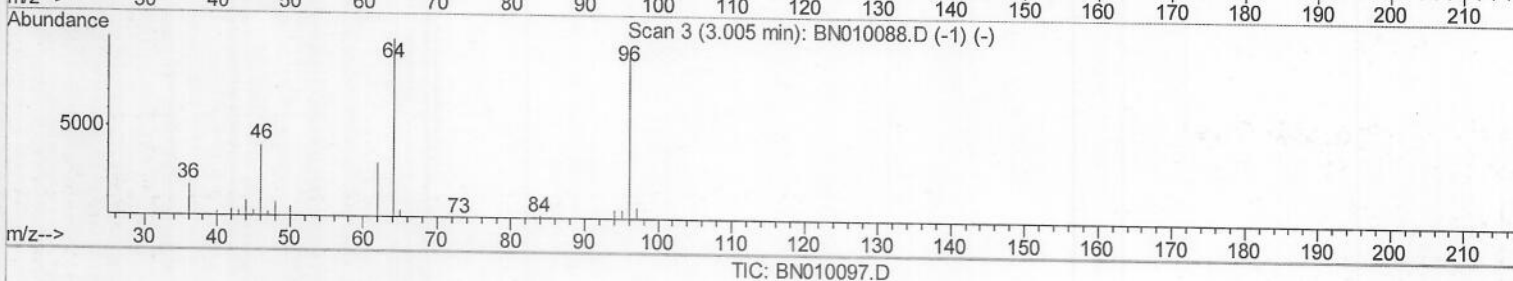
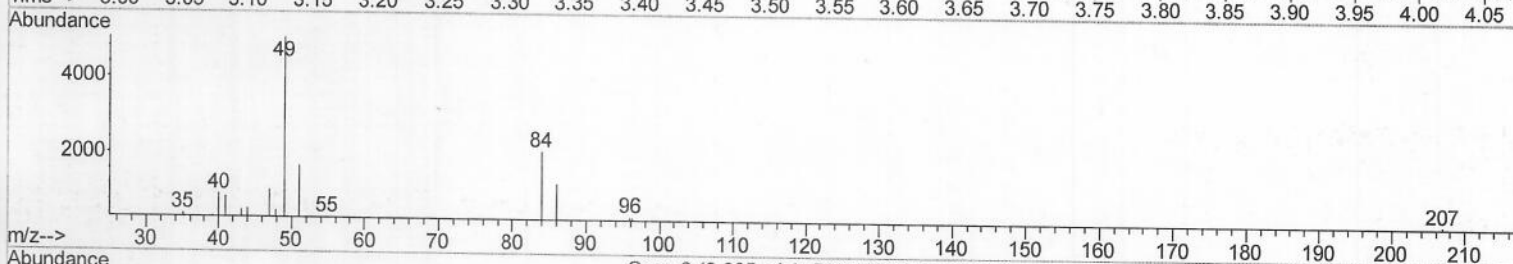
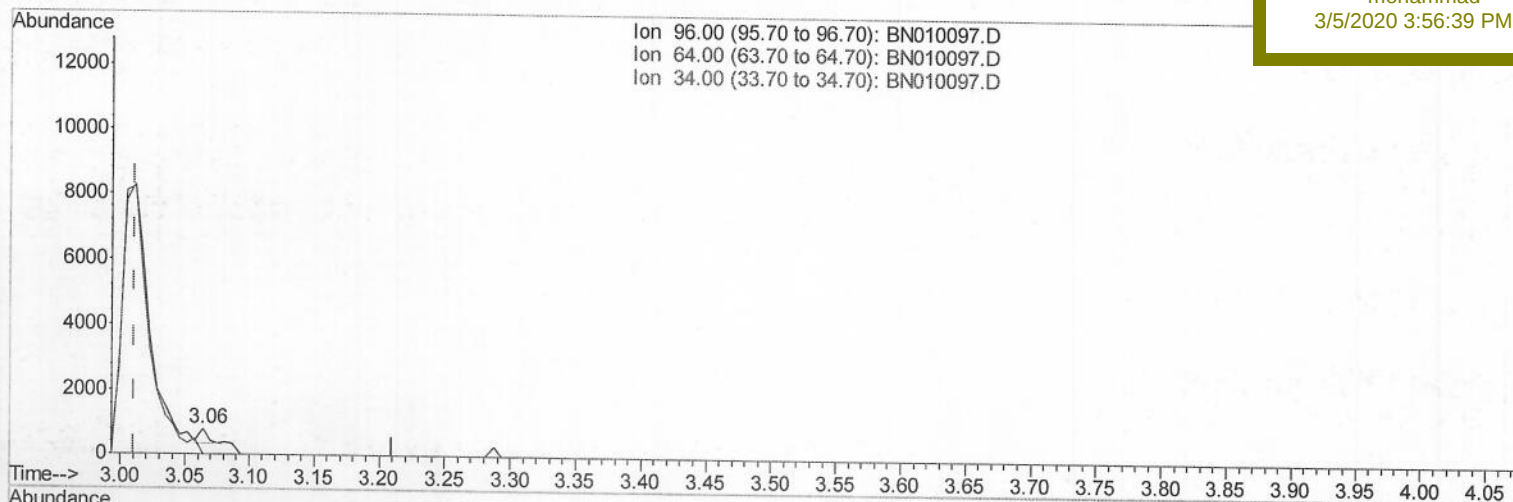
DBDS5

Manual Integrations

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TIC: BN010097.D

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

3.064min (+0.053) 0.09ng/uL

response 179

Ion	Exp%	Act%
96.00	100	100
64.00	93.60	0.00#
34.00	0.00	0.00
0.00	0.00	0.00

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BNA_N

ClientSampleId :

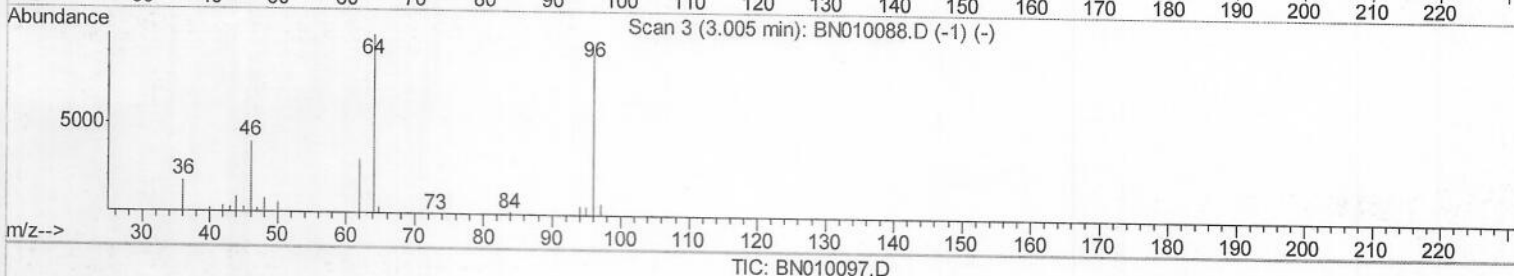
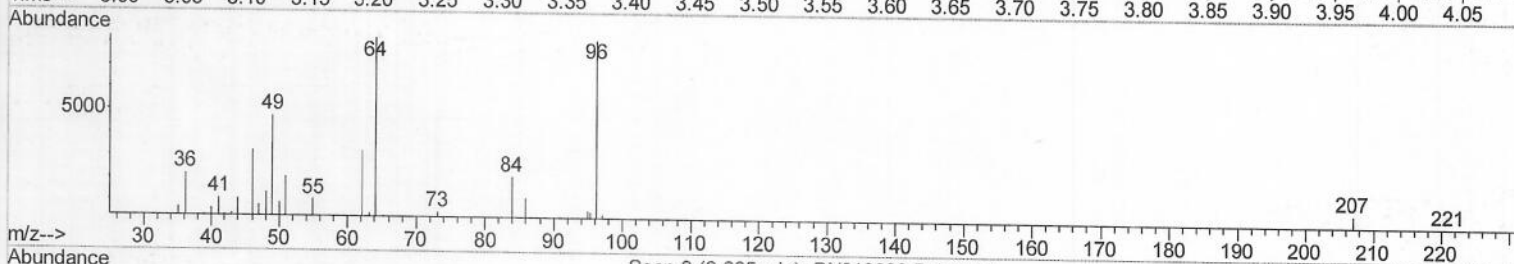
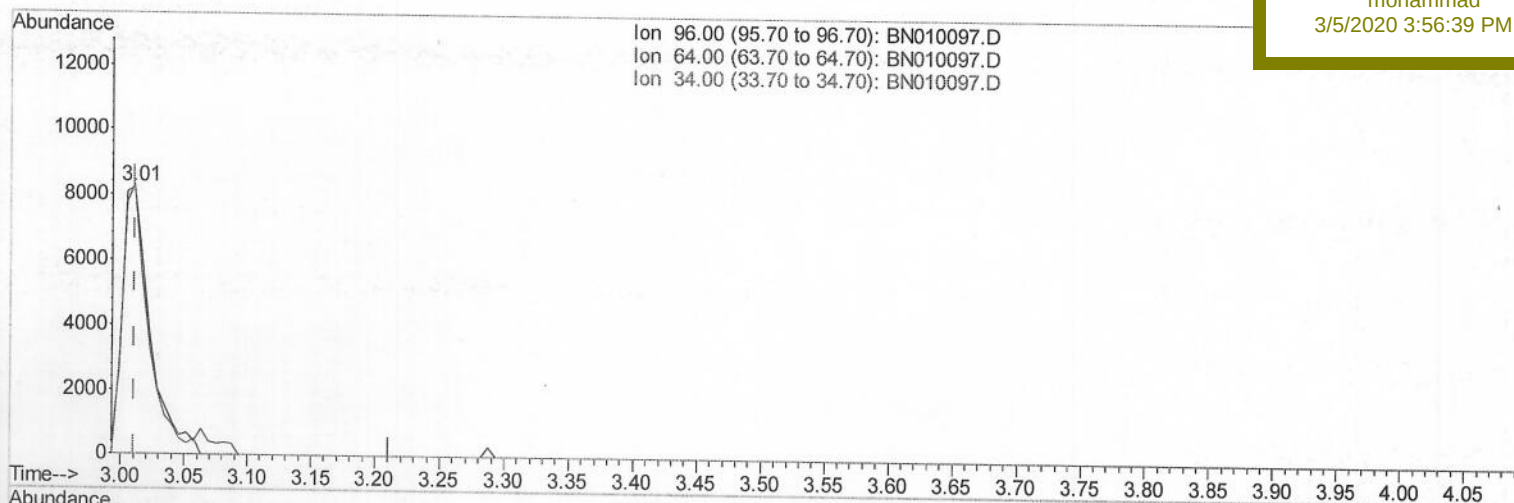
DBDS5

Manual Integrations

APPROVED

mohammad

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

3.011min (+0.000) 6.27ng/uL m

>JU 03/06/20

response 13172

Ion	Exp%	Act%
96.00	100	100
64.00	93.60	100.69
34.00	0.00	0.00
0.00	0.00	0.00

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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 04 01:30:16 2020
 Response via : Initial Calibration

Instrument :
 BNA_N
 Client Sampled :
 DBDS5

Manual Integrations
 APPROVED

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 3/5/2020 3:56:39 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.36	152	86422	20.00	ng/ul	0.00
18) Naphthalene-d8	10.10	136	373457	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.00	164	241063	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.76	188	485235	20.00	ng/ul	0.00
78) Chrysene-d12	20.99	240	415776	20.00	ng/ul	0.00
86) Perylene-d12	23.08	264	452543	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.01	96	13172m	6.27	ng/uL	0.00
5) Phenol-d5	6.56	99	264305	35.95	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.72	67	187434	36.95	ng/ul	0.00
9) 2-Chlorophenol-d4	6.90	132	210128	37.97	ng/ul	0.00
13) 4-Methylphenol-d8	8.08	113	215717	36.89	ng/ul	0.00
19) Nitrobenzene-d5	8.50	128	106237	38.88	ng/ul	0.00
22) 2-Nitrophenol-d4	9.21	143	123122	41.19	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.75	165	207930	36.19	ng/ul	0.00
29) 4-Chloroaniline-d4	10.26	131	277826	43.03	ng/ul	0.00
44) Dimethylphthalate-d6	13.43	166	678640	37.69	ng/ul	0.00
47) Acenaphthylene-d8	13.69	160	809499	38.17	ng/ul	0.00
52) 4-Nitrophenol-d4	14.26	143	100444	30.54	ng/ul	0.00
58) Fluorene-d10	15.01	176	596681	38.38	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.17	200	94695	31.41	ng/ul	0.00
71) Anthracene-d10	16.86	188	882391	39.67	ng/ul	0.00
79) Pyrene-d10	19.17	212	925439	42.59	ng/ul	0.00
90) Benzo(a)pyrene-d12	22.95	264	910312	40.28	ng/ul	0.00

25009/06/20

Target Compounds

					Ovalue	
6) Phenol	6.59	94	19778	2.511 ng/ul#	79	
45) Dimethylphthalate	13.47	163	62308	3.316 ng/ul	99	

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