

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062619\
Data File : BN006585.D
Acq On : 26 Jun 2019 19:20
Operator : HP/JU
Sample : K3498-15
Misc :
ALS Vial : 4 Sample Multiplier: 1

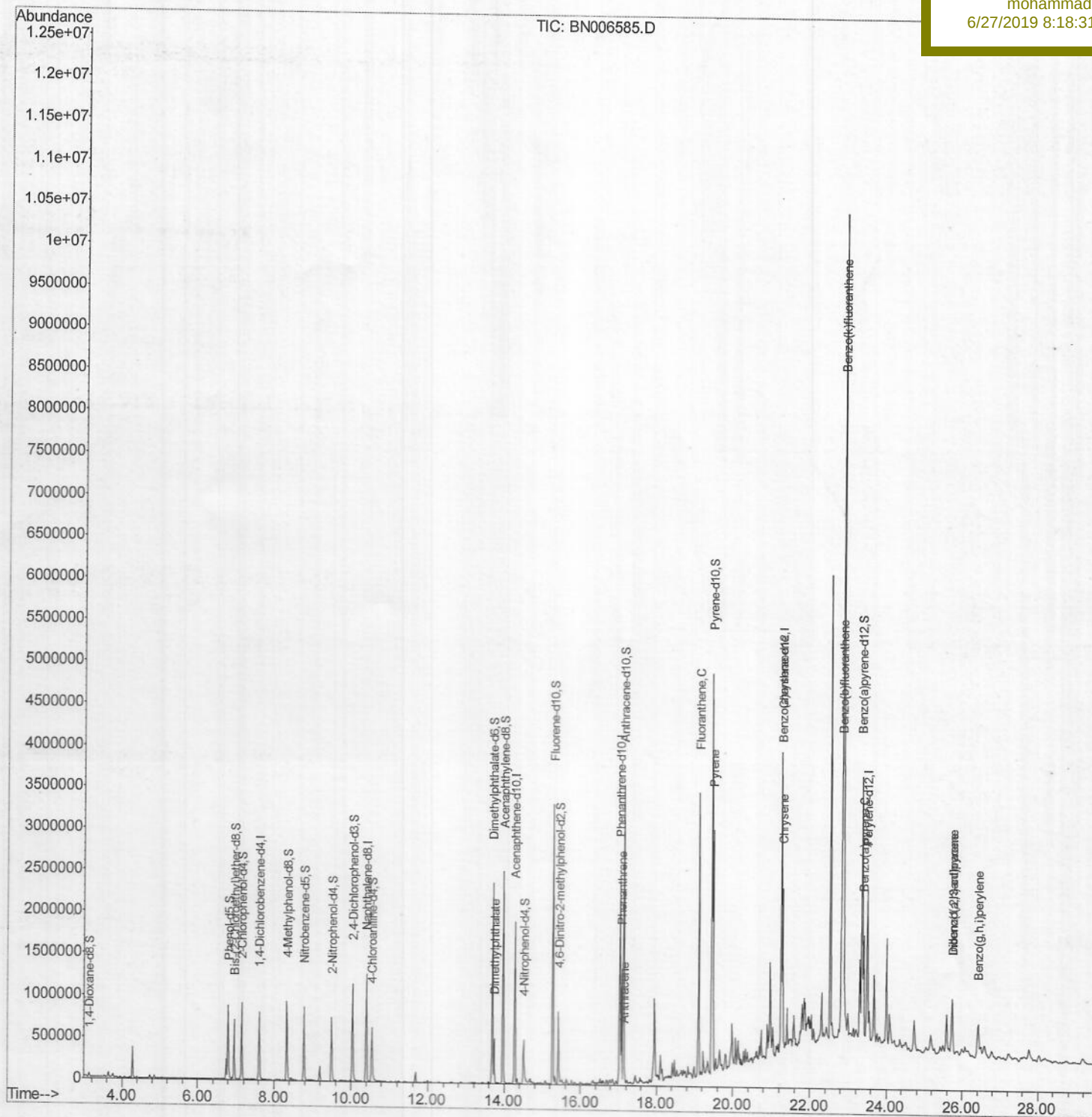
Quant Time: Jun 27 03:18:20 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 27 02:30:52 2019
Response via : Initial Calibration

Instrument :

BNA_N

Client Sampled :

C5AB4

Manual Integrations
APPROVEDmohammad
6/27/2019 8:18:31 AM

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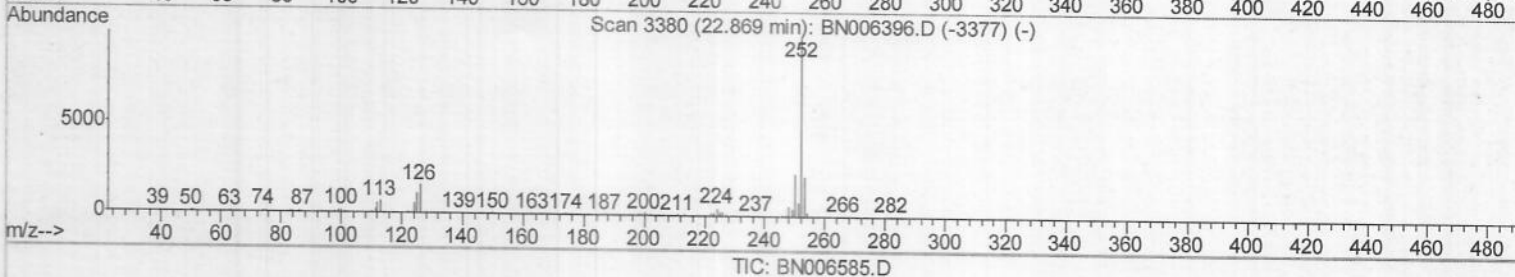
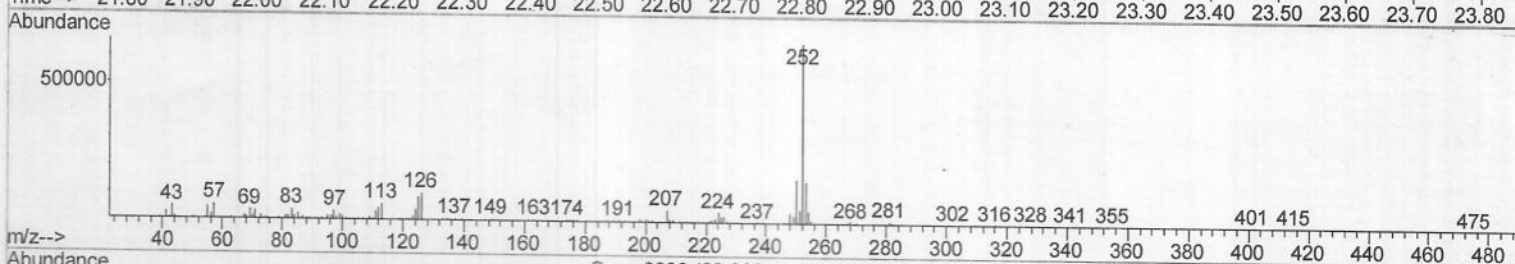
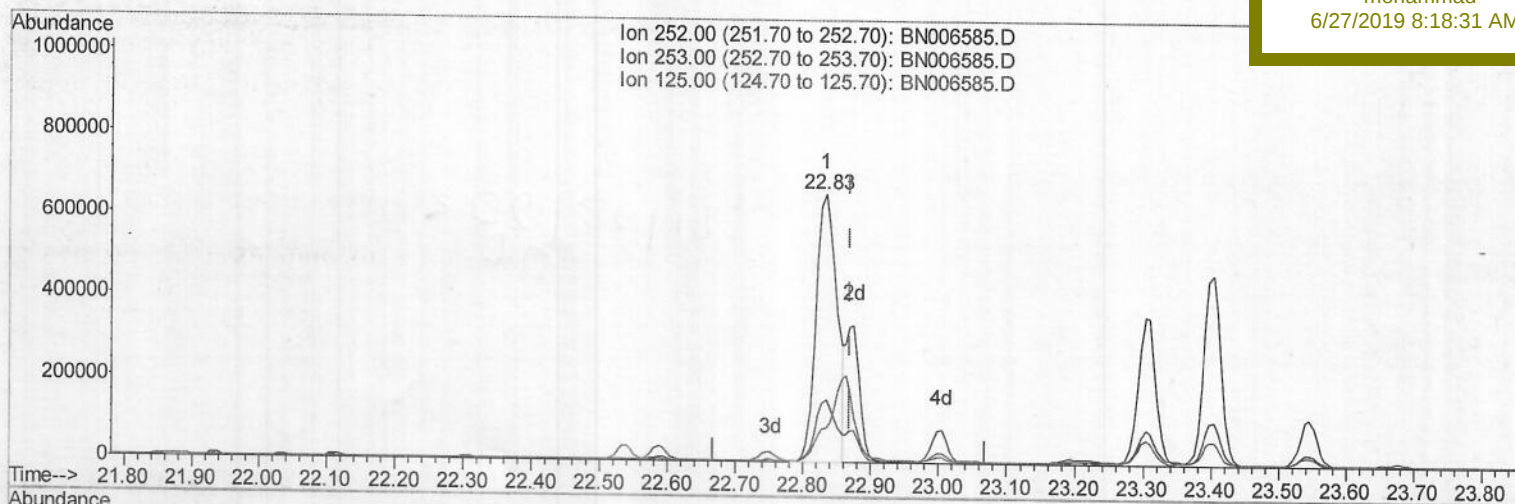
C5AB4

Manual Integrations

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(88) Benzo(k)fluoranthene

22.833min (-0.035) 19.57ng/ul

response 1387319

Ion	Exp%	Act%
252.00	100	100
253.00	21.70	23.06
125.00	11.40	12.95
0.00	0.00	0.00

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ClientSampleId :

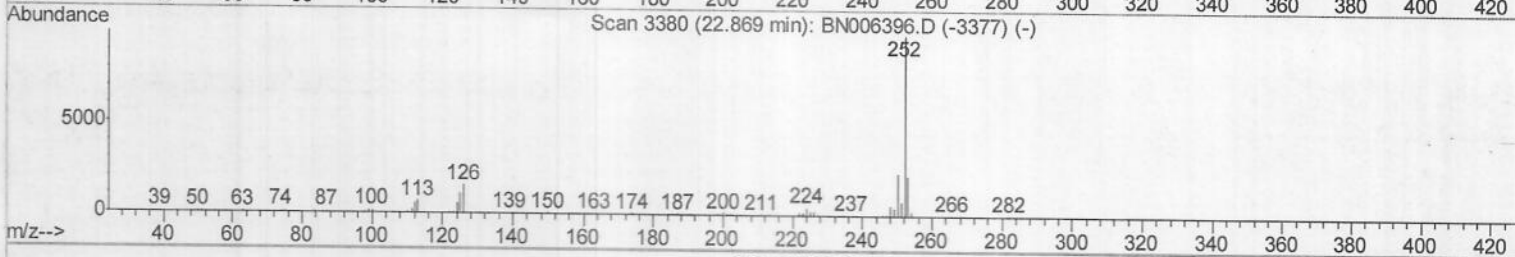
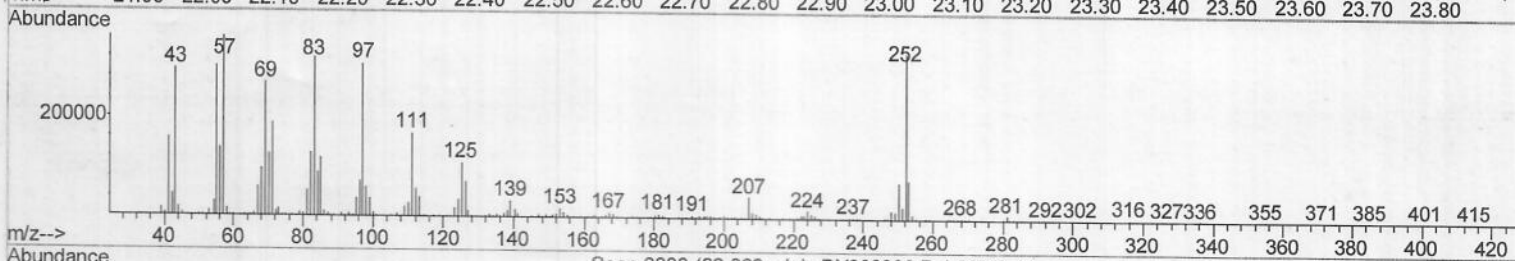
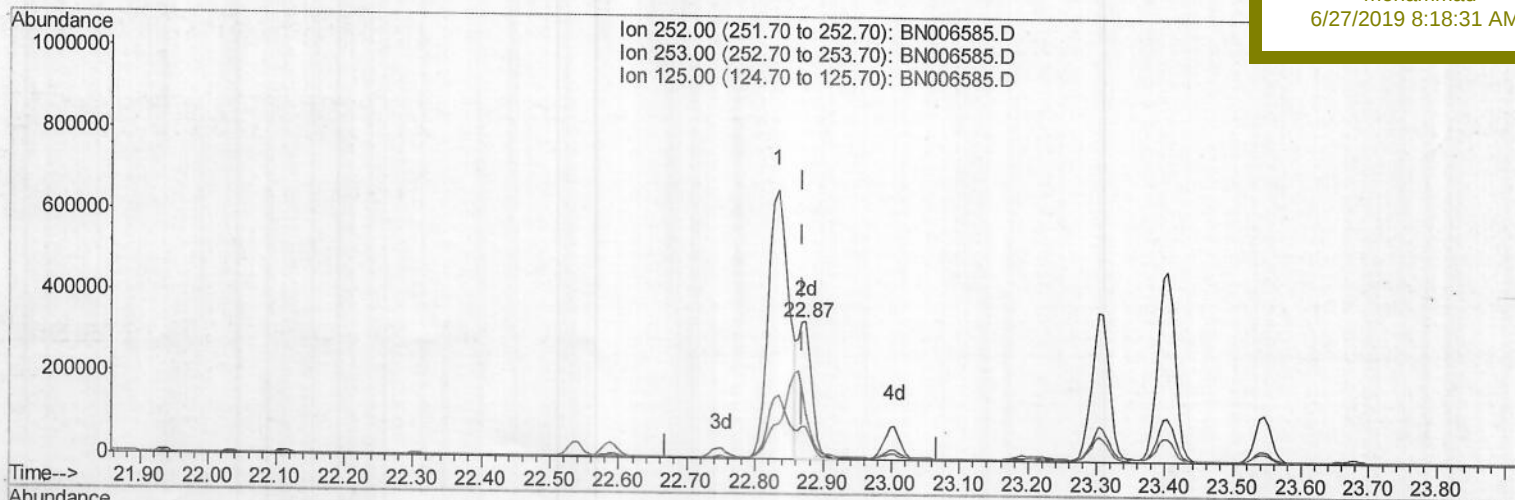
C5AB4

Manual Integrations

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TIC: BN006585.D

(88) Benzo(k)fluoranthene

22.874min (+0.006) 6.89ng/ul m

response 488835

Ion Exp% Act%

252.00 100 100

253.00 21.70 23.75

125.00 11.40 32.82#

0.00 0.00 0.00

> JU 06/27/19

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 QLast Update : Thu Jun 27 02:30:52 2019
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 BNA_N
 Client Sampled :
 C5AB4

Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	213151	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.40	136	1018245	20.00	ng/ul	-0.01
35) Acenaphthene-d10	14.27	164	602639	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.03	188	1339311	20.00	ng/ul	0.00
77) Chrysene-d12	21.26	240	1239623	20.00	ng/ul	0.00
85) Perylene-d12	23.50	264	1236049	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	22296	4.09	ng/uL	-0.01
5) Phenol-d5	6.80	99	528284	23.28	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.95	67	326844	23.26	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.15	132	404558	24.17	ng/ul	-0.01
13) 4-Methylphenol-d8	8.33	113	359297	19.96	ng/ul	-0.01
19) Nitrobenzene-d5	8.78	128	209537	25.70	ng/ul	0.00
22) 2-Nitrophenol-d4	9.49	143	240076	27.02	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	440721	25.83	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.55	131	387284	19.73	ng/ul	0.00
43) Dimethylphthalate-d6	13.69	166	1386877	26.26	ng/ul	-0.01
46) Acenaphthylene-d8	13.96	160	1670114	25.80	ng/ul	-0.01
51) 4-Nitrophenol-d4	14.51	143	164059	18.18	ng/ul	0.00
57) Fluorene-d10	15.27	176	1218217	27.12	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.42	200	165975	19.58	ng/ul	0.00
70) Anthracene-d10	17.13	188	1887392	27.99	ng/ul	0.00
78) Pyrene-d10	19.45	212	2090984	28.88	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.36	264	1954409	27.99	ng/ul	0.00

Target Compounds

						Qvalue
44) Dimethylphthalate	13.73	163	304603	6.256	ng/ul	99
69) Phenanthrene	17.07	178	758500	10.268	ng/ul	100
71) Anthracene	17.16	178	95190	1.269	ng/ul	99
76) Fluoranthene	19.10	202	1929850	24.240	ng/ul	99
79) Pyrene	19.47	202	1594541	18.893	ng/ul	100
82) Benzo(a)anthracene	21.25	228	861109	10.163	ng/ul	97
84) Chrysene	21.30	228	1091654	13.538	ng/ul	99
87) Benzo(b)fluoranthene	22.83	252	1387319	18.701	ng/ul	97
88) Benzo(k)fluoranthene	22.87	252	488835m	6.895	ng/ul	
90) Benzo(a)pyrene	23.40	252	823433	11.576	ng/ul	100
91) Indeno(1,2,3-cd)pyrene	25.73	276	550935	6.545	ng/ul	95
92) Dibenzo(a,h)anthracene	25.73	278	140882	2.021	ng/ul#	78
93) Benzo(g,h,i)perylene	26.42	276	457584	6.454	ng/ul	99

JU 06/27/19

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