

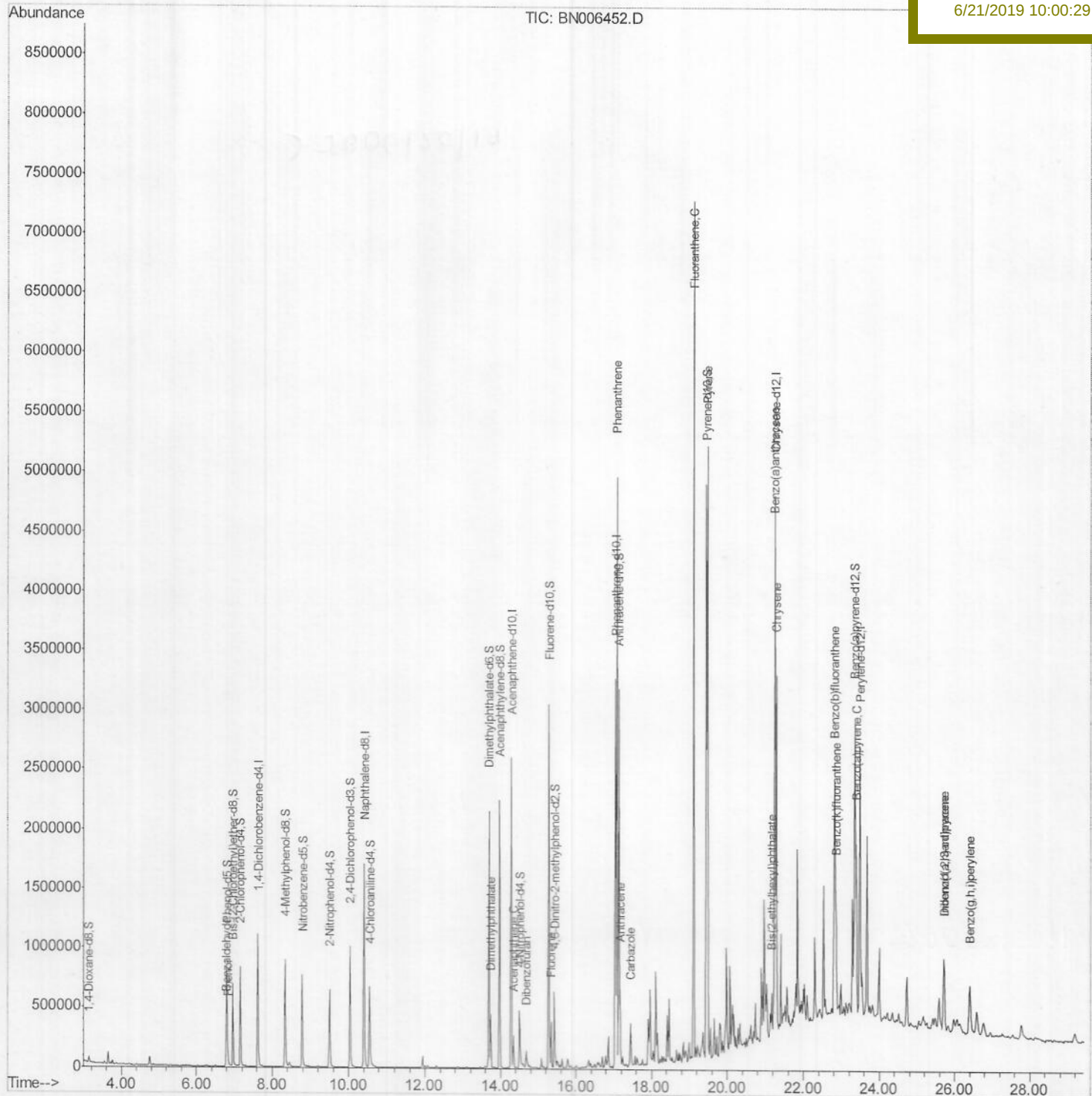
Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN061919\
Data File : BN006452.D
Acq On : 21 Jun 2019 05:46
Operator : JU/SJ
Sample : K3360-14
Misc :
ALS Vial : 57 Sample Multiplier: 1

Quant Time: Jun 21 06:41:54 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 21 06:31:50 2019
Response via : Initial Calibration

Instrument :
BNA_N
Client Sample ID :
A41T6

Manual Integrations
APPROVED

mohammad
6/21/2019 10:00:29 PM



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Operator : JU/SJ

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QLast Update : Fri Jun 21 06:31:50 2019

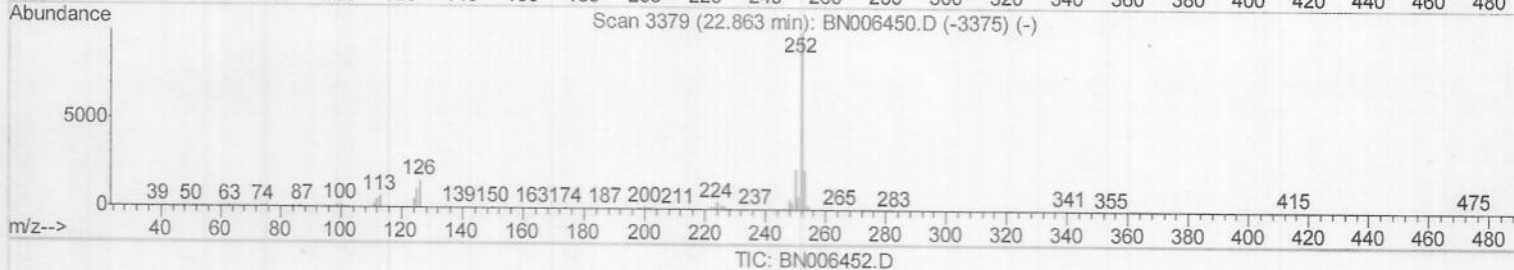
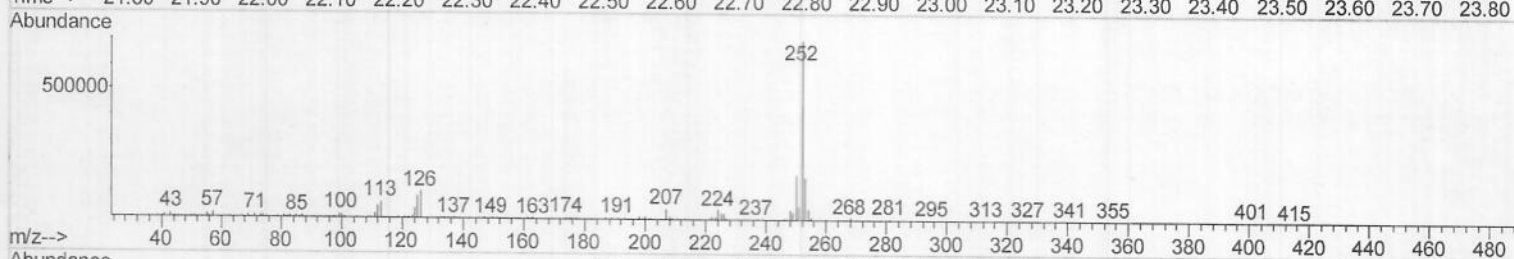
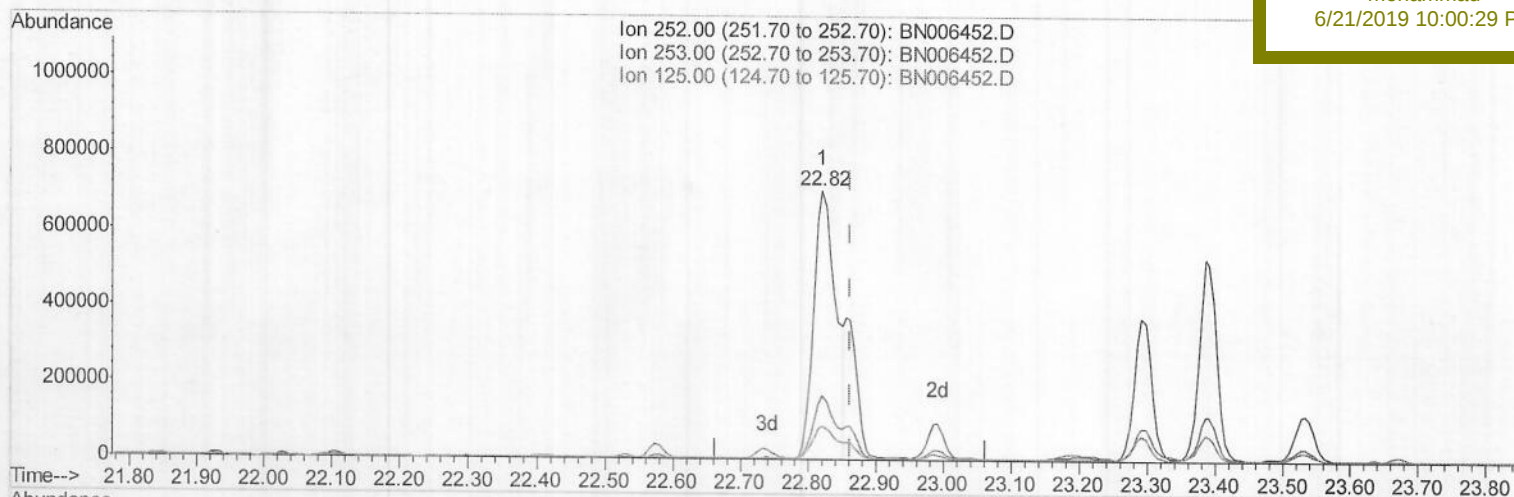
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(88) Benzo(k)fluoranthene

22.821min (-0.042) 18.44ng/ul

response 1580221

Ion Exp% Act%

252.00 100 100

253.00 21.70 23.52

125.00 11.40 12.42

0.00 0.00 0.00

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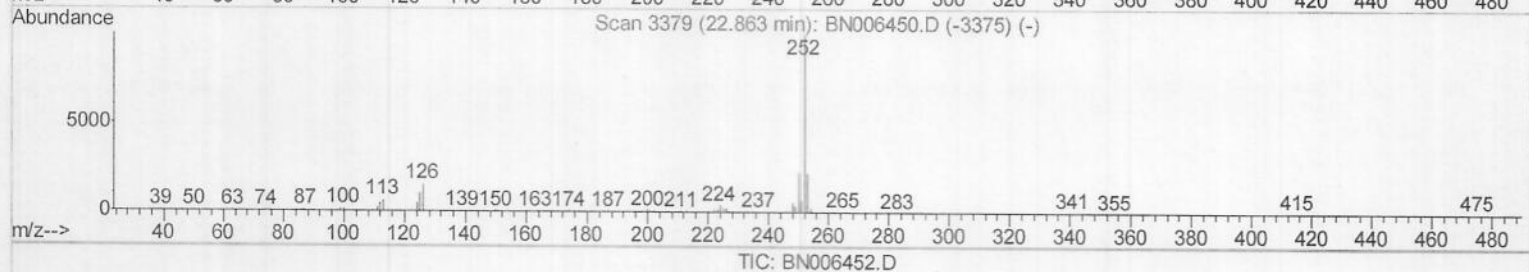
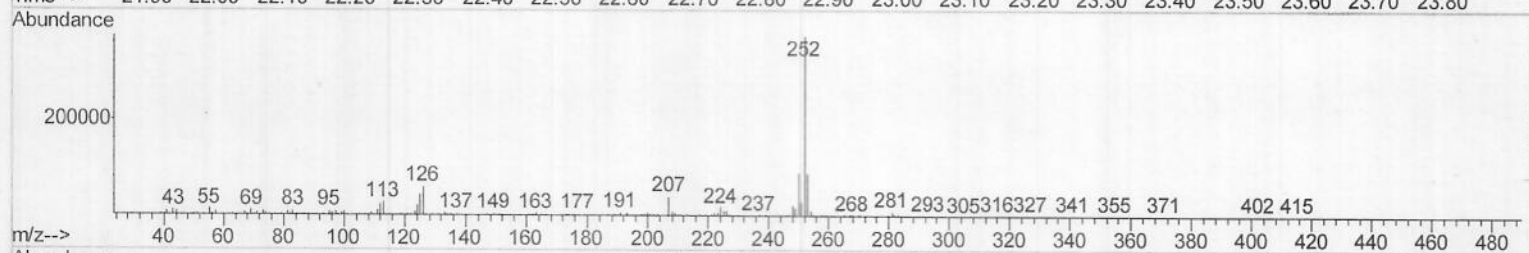
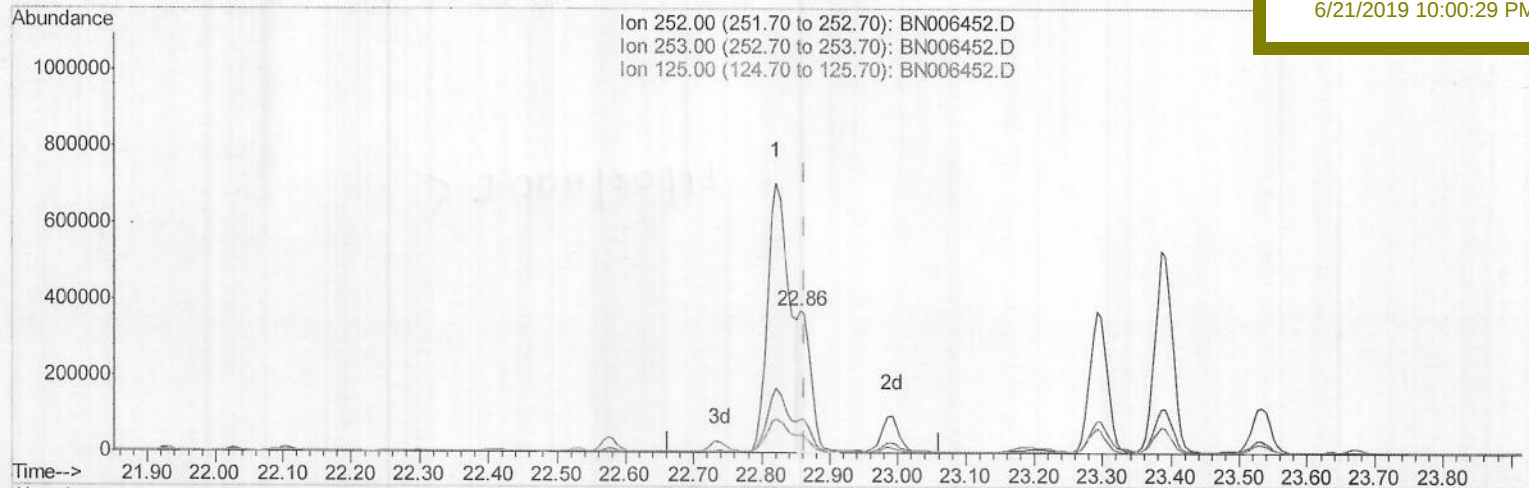
Response via : Initial Calibration

Instrument :

BNA_N

ClientSampleId :

A41T6

Manual Integrations
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6/21/2019 10:00:29 PM

(88) Benzo(k)fluoranthene

22.856min (-0.006) 5.57ng/ul m) JU06120119

response 477421

Ion	Exp%	Act%
252.00	100	100
253.00	21.70	23.18
125.00	11.40	11.76
0.00	0.00	0.00

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 Sample : K3360-14
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_N
 Client Sampled :
 A41T6

Manual Integrations
 APPROVED

mohammad
 6/21/2019 10:00:29 PM

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 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 21 06:31:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	283339	20.00	ng/ul	0.00
18) Naphthalene-d8	10.40	136	1334147	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.28	164	812303	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.04	188	1821668	20.00	ng/ul	0.00
77) Chrysene-d12	21.26	240	1602757	20.00	ng/ul	0.00
85) Perylene-d12	23.49	264	1494171	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	22618	3.12	ng/uL	0.00
5) Phenol-d5	6.80	99	478796	15.87	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.96	67	318361	17.05	ng/ul	0.00
9) 2-Chlorophenol-d4	7.16	132	362521	16.30	ng/ul	0.00
13) 4-Methylphenol-d8	8.33	113	339713	14.19	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	179359	16.79	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	200107	17.19	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	363671	16.26	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	385354	14.99	ng/ul	0.00
43) Dimethylphthalate-d6	13.69	166	1316240	18.49	ng/ul	0.00
46) Acenaphthylene-d8	13.97	160	1509653	17.30	ng/ul	0.00
51) 4-Nitrophenol-d4	14.50	143	136596	11.23	ng/ul	0.00
57) Fluorene-d10	15.28	176	1138588	18.80	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.42	200	114581	9.94	ng/ul	0.00
70) Anthracene-d10	17.13	188	1711869	18.67	ng/ul	0.00
78) Pyrene-d10	19.44	212	1843057	19.69	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.34	264	1547739	18.34	ng/ul	0.00

Target Compounds

					Qvalue	
4) Benzaldehyde	6.78	77	13298	1.091	ng/ul	91
6) Phenol	6.83	94	32961	1.133	ng/ul#	84
44) Dimethylphthalate	13.74	163	261605	3.986	ng/ul	99
49) Acenaphthene	14.35	153	101228	1.837	ng/ul	99
53) Dibenzofuran	14.68	168	82408	1.079	ng/ul	98
58) Fluorene	15.33	166	153632	2.489	ng/ul	99
69) Phenanthrene	17.08	178	2753711	27.407	ng/ul	100
71) Anthracene	17.17	178	387126	3.796	ng/ul	99
74) Carbazole	17.44	167	209768	2.383	ng/ul	100
76) Fluoranthene	19.11	202	4044641	37.352	ng/ul	98
79) Pyrene	19.47	202	3121878	28.609	ng/ul	99
82) Benzo(a)anthracene	21.24	228	1452961	13.263	ng/ul	98
83) Bis(2-ethylhexyl)phthalate	21.17	149	92342	1.316	ng/ul	99
84) Chrysene	21.29	228	1542584	14.795	ng/ul	98
87) Benzo(b)fluoranthene	22.82	252	1580221	17.622	ng/ul	97
88) Benzo(k)fluoranthene	22.86	252	477421m)	5.571	ng/ul	100
90) Benzo(a)pyrene	23.39	252	940837	10.942	ng/ul	95
91) Indeno(1,2,3-cd)pyrene	25.72	276	543924	1.732	ng/ul#	80
92) Dibenzo(a,h)anthracene	25.72	278	145993	5.412	ng/ul	98
93) Benzo(g,h,i)perylene	26.40	276	463842			

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