

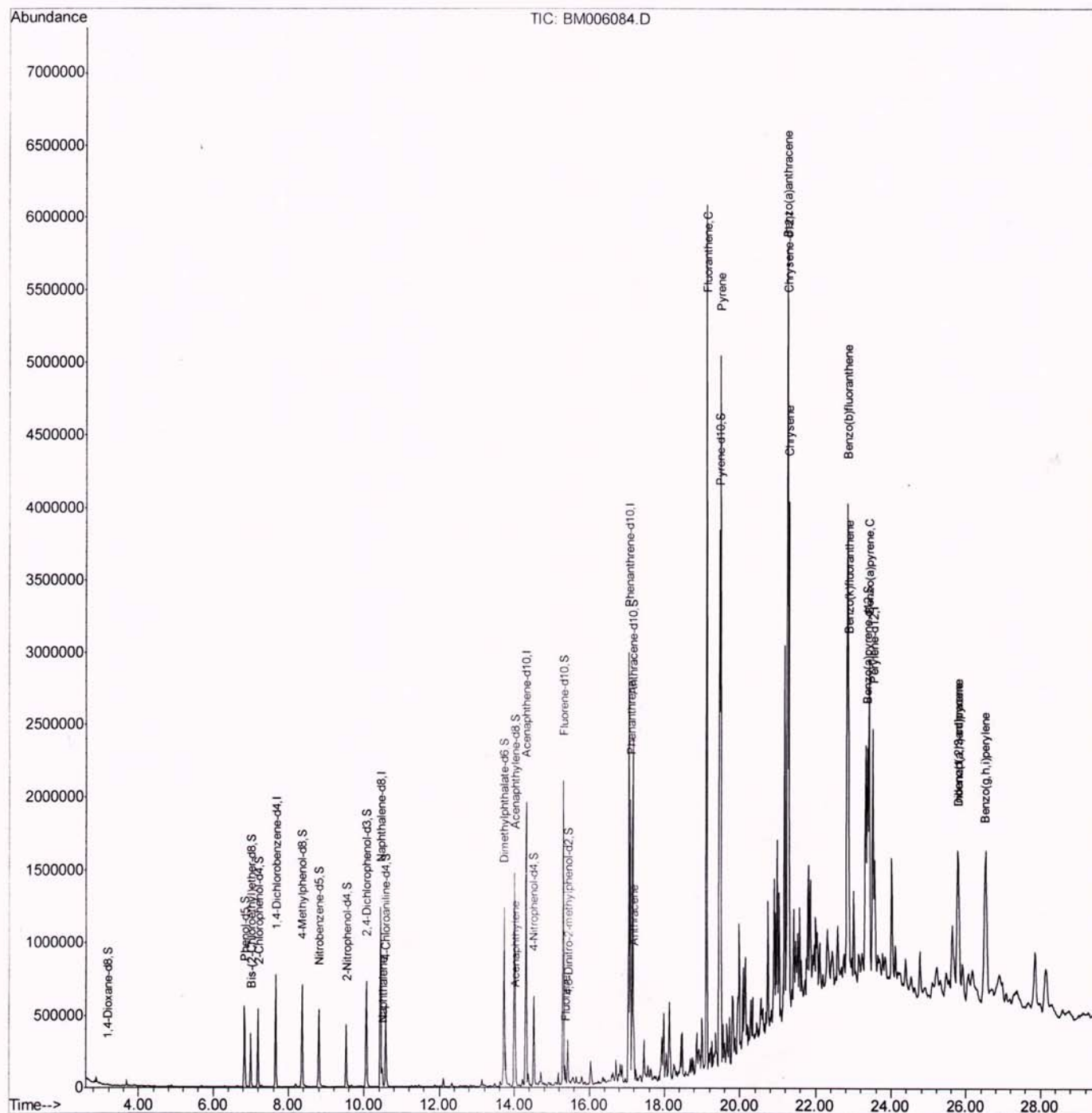
Instrument :
BNA_M
ClientSampleId :
D9YD4

Manual Integrations APPROVED

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```
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062716\  
Data File : BM006084.D  
Acq On    : 27 Jun 2016   23:45  
Operator  : UM/SJ  
Sample    : H3779-03  
Misc      :  
ALS Vial  : 18      Sample Multiplier: 1
```

Quant Time: Jun 28 04:04:01 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 28 02:11:41 2016
Response via : Initial Calibration



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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	198922	20.00	ng/ul	0.00
7) Naphthalene-d8	10.44	136	1005286	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.31	164	660543	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.06	188	1676396	20.00	ng/ul	0.00
29) Chrysene-d12	21.26	240	1746841	20.00	ng/ul	0.00
34) Perylene-d12	23.51	264	1645497	20.00	ng/ul	0.03

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.18	96	5448	1.22	ng/uL	0.00
3) Phenol-d5	6.83	99	305720	14.61	ng/ul	0.00
4) Bis-(2-Chloroethyl) ether-d	7.00	67	174303	14.97	ng/ul	0.00
5) 2-Chlorophenol-d4	7.19	132	226552	15.54	ng/ul	0.00
6) 4-Methylphenol-d8	8.36	113	257673	14.59	ng/ul	0.00
8) Nitrobenzene-d5	8.81	128	115437	15.08	ng/ul	0.00
9) 2-Nitrophenol-d4	9.53	143	124807	14.66	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.06	165	258133	15.80	ng/ul	0.00
12) 4-Chloroaniline-d4	10.57	131	290020	16.36	ng/ul	0.00
16) Dimethylphthalate-d6	13.72	166	820492	14.51	ng/ul	0.00
17) Acenaphthylene-d8	14.00	160	1076734	16.42	ng/ul	0.00
20) 4-Nitrophenol-d4	14.51	143	136458	11.31	ng/ul	-0.02
21) Fluorene-d10	15.30	176	780353	16.16	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.43	200	60669	6.52	ng/ul	-0.01
26) Anthracene-d10	17.16	188	1299910	16.88	ng/ul	0.00
30) Pyrene-d10	19.46	212	1488016	19.45	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.36	264	1220600	16.43	ng/ul	0.02

Target Compounds

Qvalue

11) Naphthalene	10.49	128	96626	1.92	ng/ul	99
18) Acenaphthylene	14.03	152	233563	3.38	ng/ul	99
22) Fluorene	15.36	166	54396	1.03	ng/ul	99
25) Phenanthrene	17.10	178	1158230	12.84	ng/ul	100
27) Anthracene	17.19	178	365477	3.93	ng/ul	99
28) Fluoranthene	19.12	202	3692883	31.97	ng/ul	98
31) Pyrene	19.49	202	3084444	30.84	ng/ul	96
32) Benzo(a)anthracene	21.24	228	2102932	20.27	ng/ul	96
33) Chrysene	21.29	228	2028654	21.40	ng/ul	97
35) Benzo(b)fluoranthene	22.83	252	2973390	31.10	ng/ul	98
36) Benzo(k)fluoranthene	22.86	252	1011892m	11.38	ng/ul	97
38) Benzo(a)pyrene	23.40	252	1891046	20.63	ng/ul	97
39) Indeno(1,2,3-cd)pyrene	25.78	276	1437940	13.16	ng/ul#	94
40) Dibenzo(a,h)anthracene	25.77	278	358310	3.94	ng/ul	94
41) Benzo(g,h,i)perylene	26.50	276	1203906	12.74	ng/ul	98

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

U.M.
 06/28/16

Quantitation Report (Qedit)

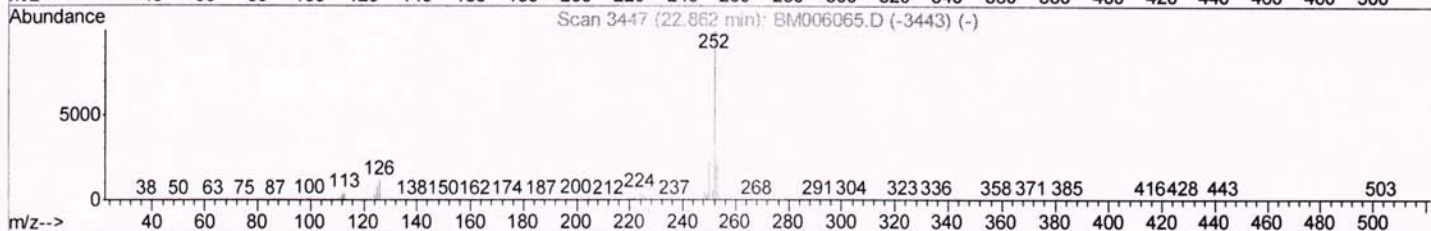
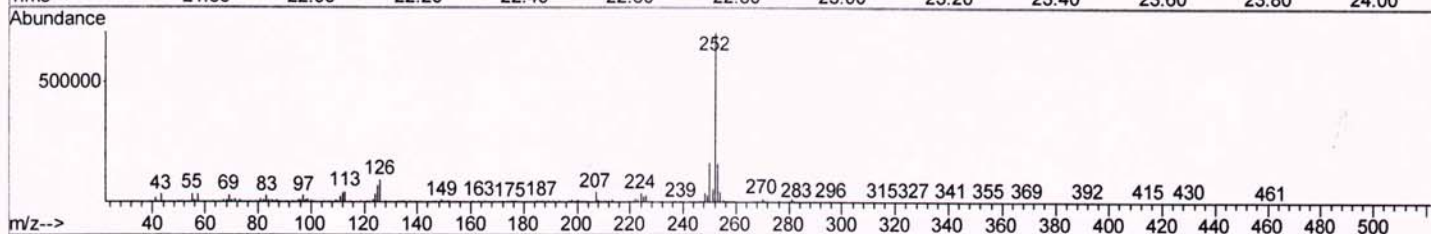
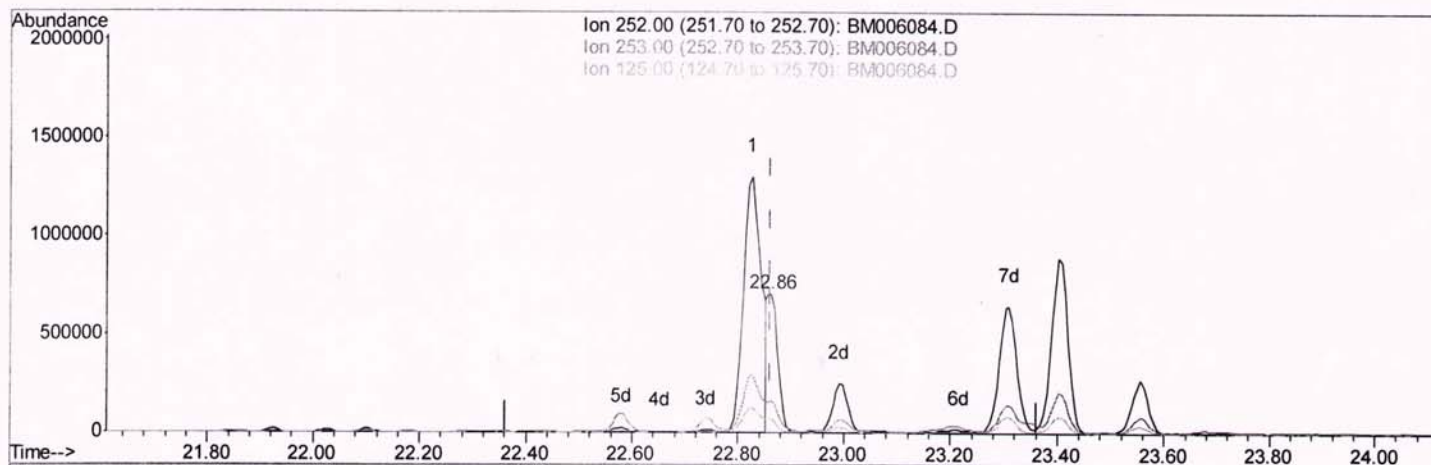
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TIC: BM006084.D

(36) Benzo(k)fluoranthene

22.862min (+0.000) 11.38ng/ul m

response 1011892

Ion Exp% Act%

252.00 100 100

253.00 22.00 22.87

125.00 10.10 9.80

0.00 0.00 0.00

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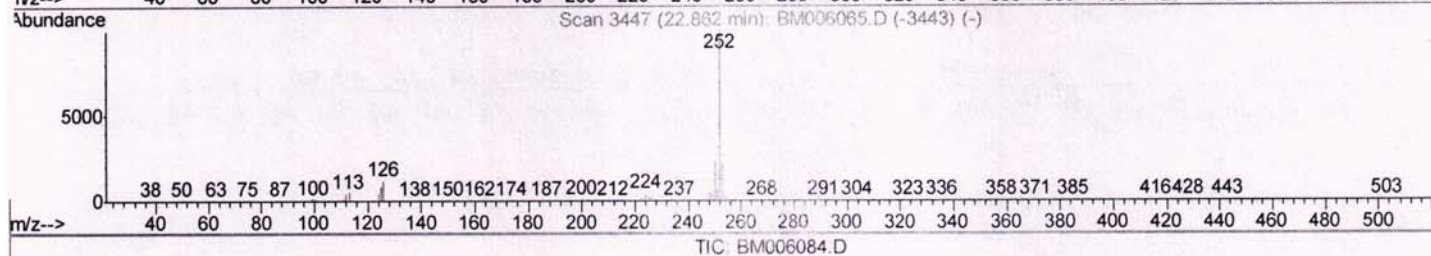
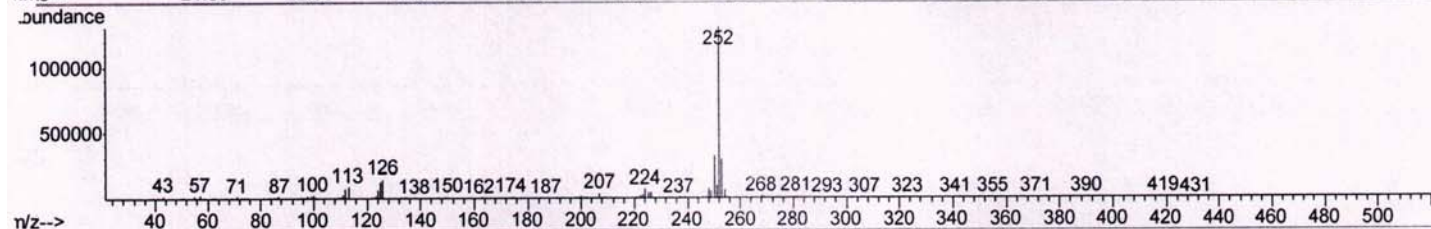
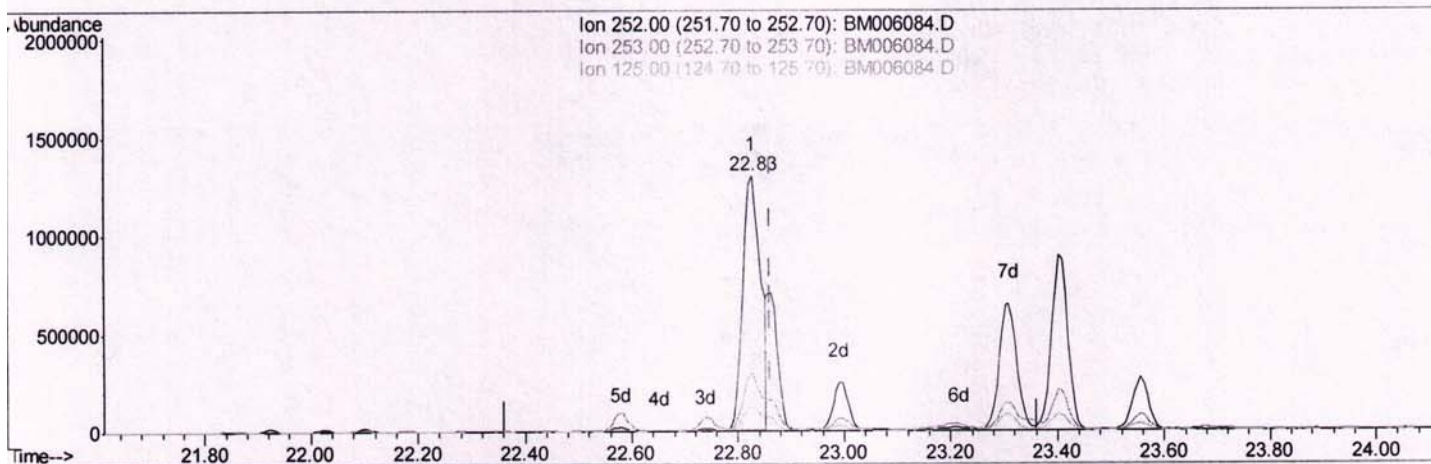
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(36) Benzo(k)fluoranthene

22.827min (-0.035) 33.43ng/ul

response 2973390

Ion	Exp%	Act%
252.00	100	100
253.00	22.00	22.88
125.00	10.10	9.69
0.00	0.00	0.00