

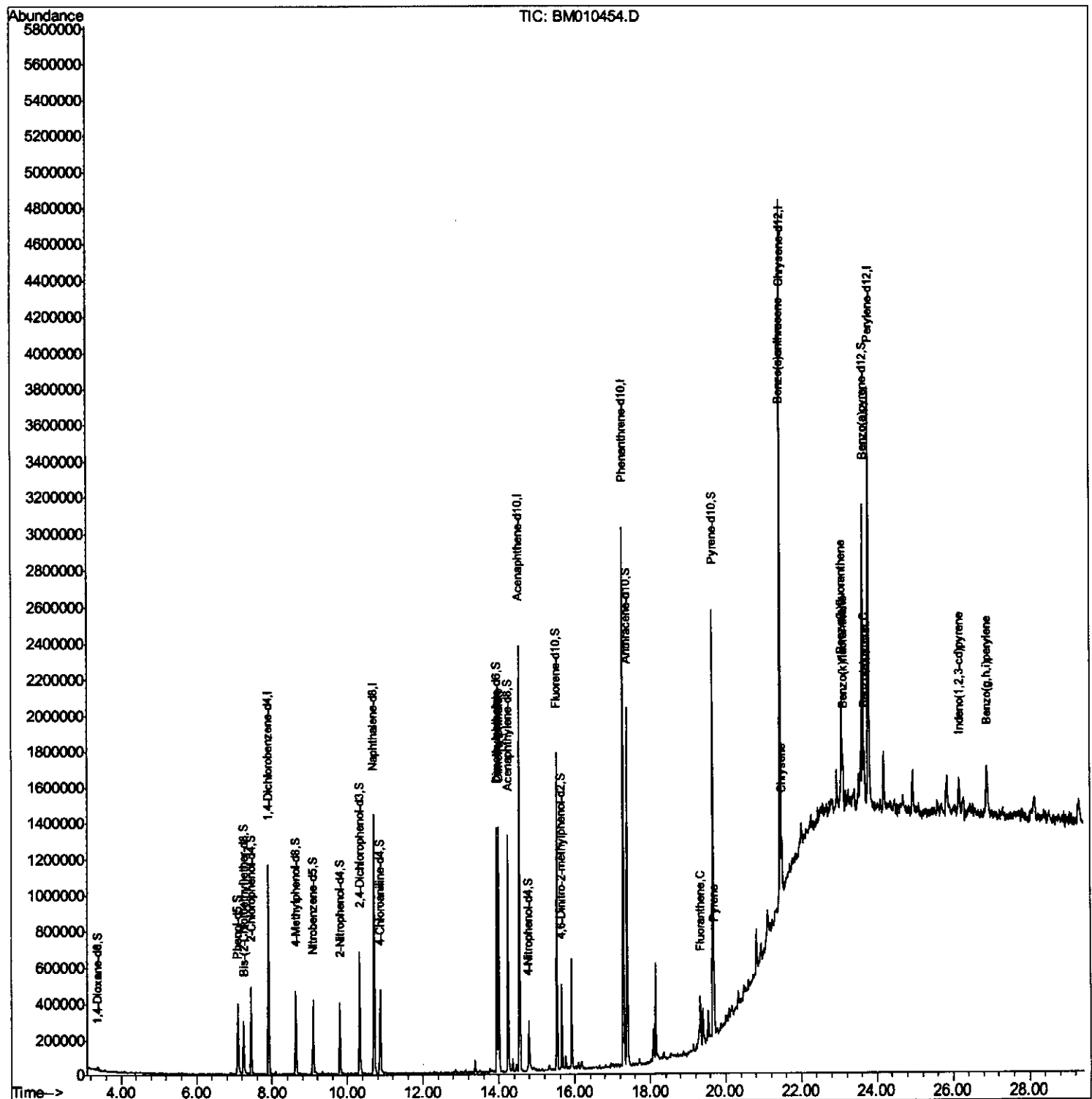
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060917\
Data File : BM010454.D
Acq On : 09 Jun 2017 15:27
Operator : SJ/MA
Sample : I3521-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F5C42

Manual Integrations
APPROVED

mohammad
6/12/2017 11:18:00 AM

Quant Time: Jun 10 02:20:58 2017
Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jun 10 01:45:58 2017
Response via : Initial Calibration



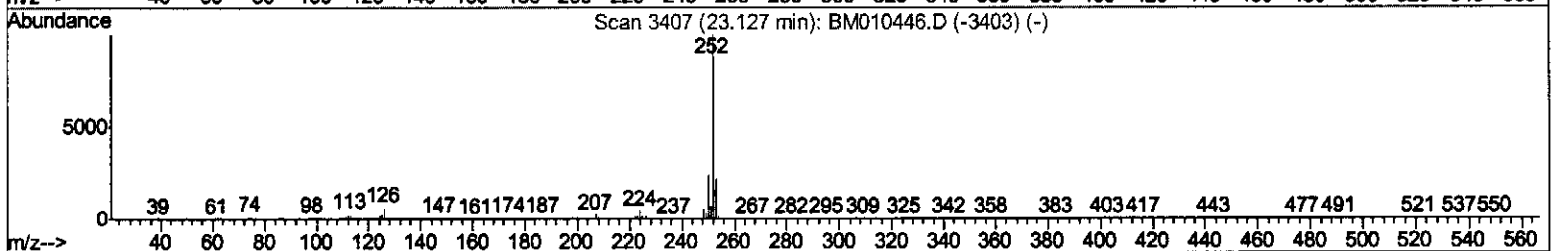
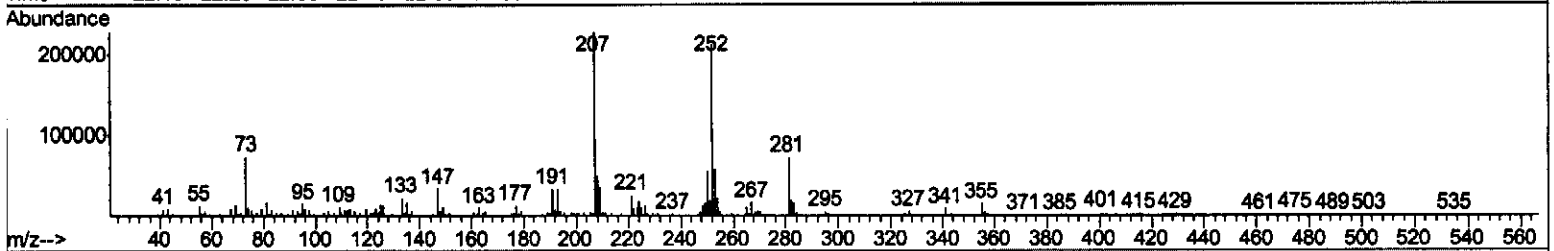
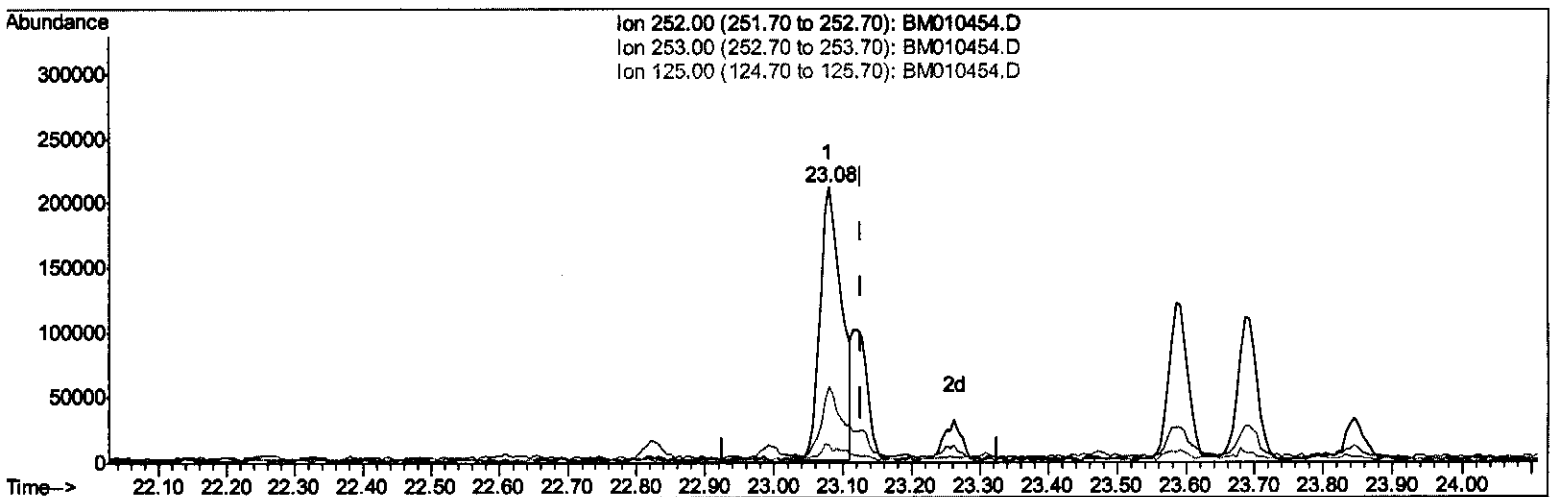
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TIC: BM010454.D

(86) Benzo(k)fluoranthene

23.080min (-0.047) 4.76ng/ul

response 457814

Ion	Exp%	Act%
252.00	100	100
253.00	21.40	27.41#
125.00	4.30	6.48#
0.00	0.00	0.00

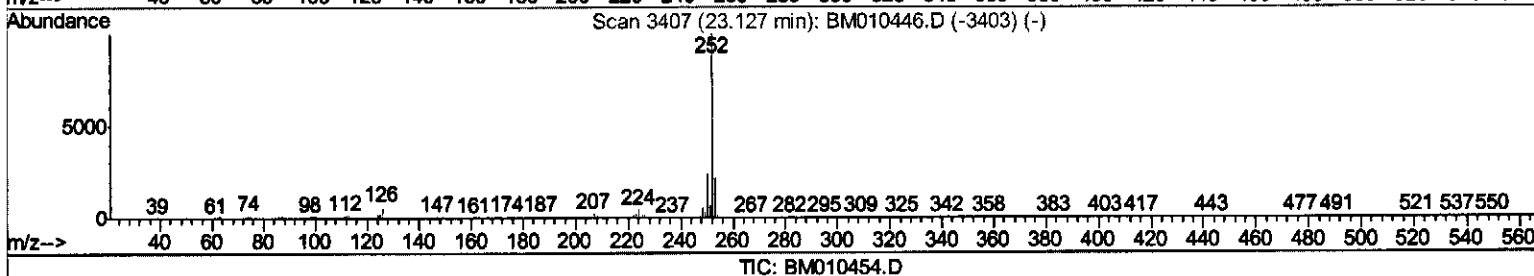
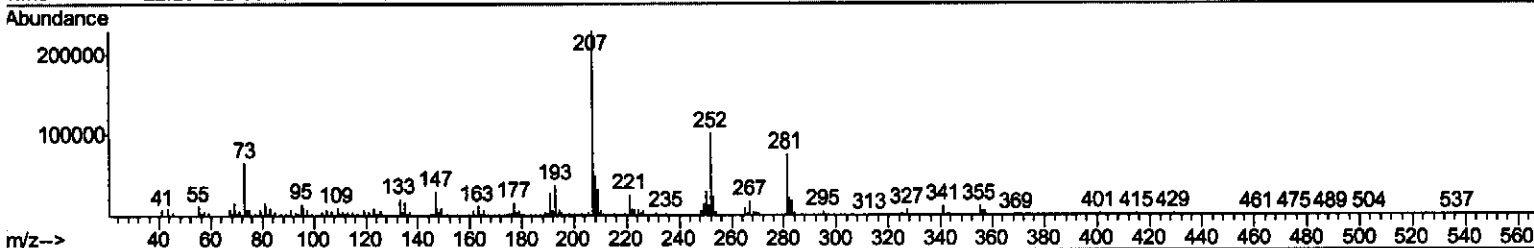
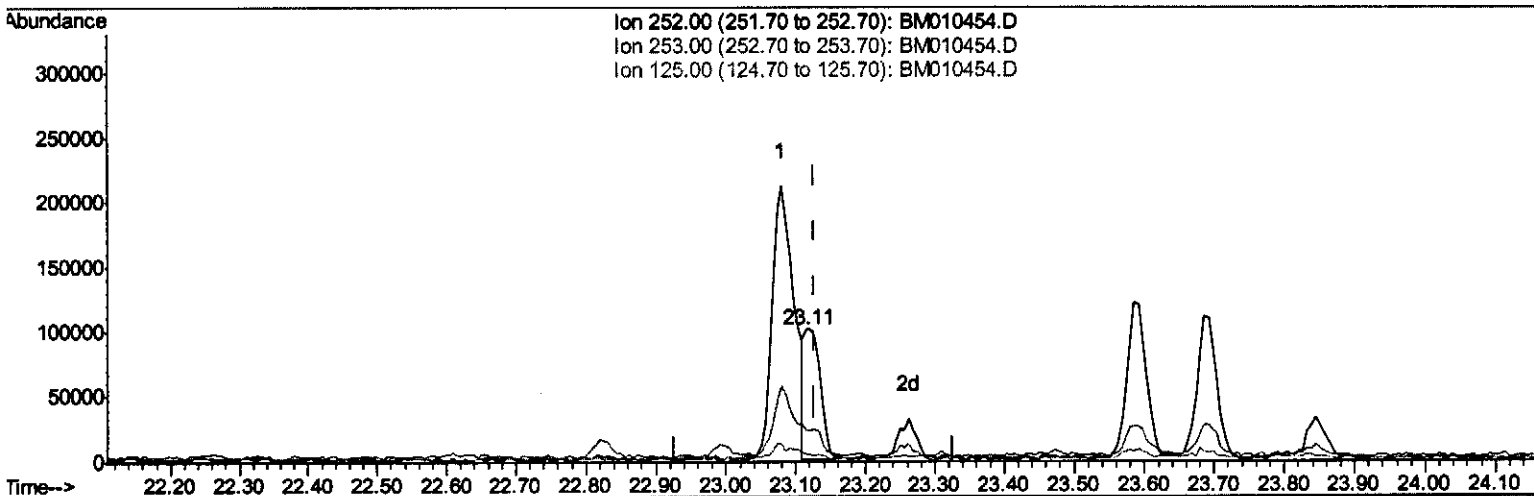
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Manual Integrations
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QLast Update : Sat Jun 10 01:45:58 2017
Response via : Initial Calibration



(86) Benzo(k)fluoranthene

23.115min (-0.012) 1.59ng/ul m

response 153284

Ion	Exp%	Act%
252.00	100	100
253.00	21.40	23.55
125.00	4.30	6.02#
0.00	0.00	0.00

SJ
06/14/17

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060917\
 Data File : BM010454.D
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 Operator : SJ/MA
 Sample : I3521-04
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F5C42

Manual Integrations
 APPROVED

mohammad
 6/12/2017 11:18:00 AM

Quant Time: Jun 10 02:20:58 2017
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 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jun 10 01:45:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	264535	20.00	ng/ul	0.00
18) Naphthalene-d8	10.71	136	997707	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.54	164	667426	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.29	188	1597047	20.00	ng/ul	0.00
75) Chrysene-d12	21.45	240	1751534	20.00	ng/ul	0.00
83) Perylene-d12	23.80	264	1730851	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.38	96	13595	2.11	ng/uL	0.00
5) Phenol-d5	7.09	99	187641	9.36	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.25	67	116521	10.34	ng/ul	0.00
9) 2-Chlorophenol-d4	7.45	132	173627	10.87	ng/ul	0.00
13) 4-Methylphenol-d8	8.63	113	148408	9.18	ng/ul	0.00
19) Nitrobenzene-d5	9.09	128	83292	11.40	ng/ul	0.00
22) 2-Nitrophenol-d4	9.80	143	100824	11.91	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.33	165	216767	12.46	ng/ul	0.00
29) 4-Chloroaniline-d4	10.87	131	193060	11.51	ng/ul	0.00
43) Dimethylphthalate-d6	13.95	166	749291	13.51	ng/ul	0.00
46) Acenaphthylene-d8	14.24	160	801980	12.40	ng/ul	0.00
51) 4-Nitrophenol-d4	14.79	143	62796	7.57	ng/ul	0.00
57) Fluorene-d10	15.53	176	635353	13.03	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.66	200	118389	10.47	ng/ul	0.00
70) Anthracene-d10	17.39	188	1014229	13.05	ng/ul	0.00
76) Pyrene-d10	19.67	212	1176046	16.24	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.64	264	1105725	13.72	ng/ul	0.00

Target Compounds

					Qvalue	
44) Dimethylphthalate	14.00	163	709075	13.002	ng/ul	99
74) Fluoranthene	19.34	202	149074	1.422	ng/ul#	94
77) Pyrene	19.70	202	195821	2.170	ng/ul#	94
80) Benzo(a)anthracene	21.44	228	159951	1.619	ng/ul	96
82) Chrysene	21.49	228	141441	1.584	ng/ul	97
85) Benzo(b)fluoranthene	23.08	252	457814	4.546	ng/ul#	90
86) Benzo(k)fluoranthene	23.11	252	153284m	1.593	ng/ul	
88) Benzo(a)pyrene	23.69	252	213768	2.141	ng/ul#	95
89) Indeno(1,2,3-cd)pyrene	26.19	276	189606	1.503	ng/ul#	97
91) Benzo(g,h,i)perylene	26.92	276	150328	1.447	ng/ul#	89

5506/19/17

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