

Quantitation Report (QT Reviewed)

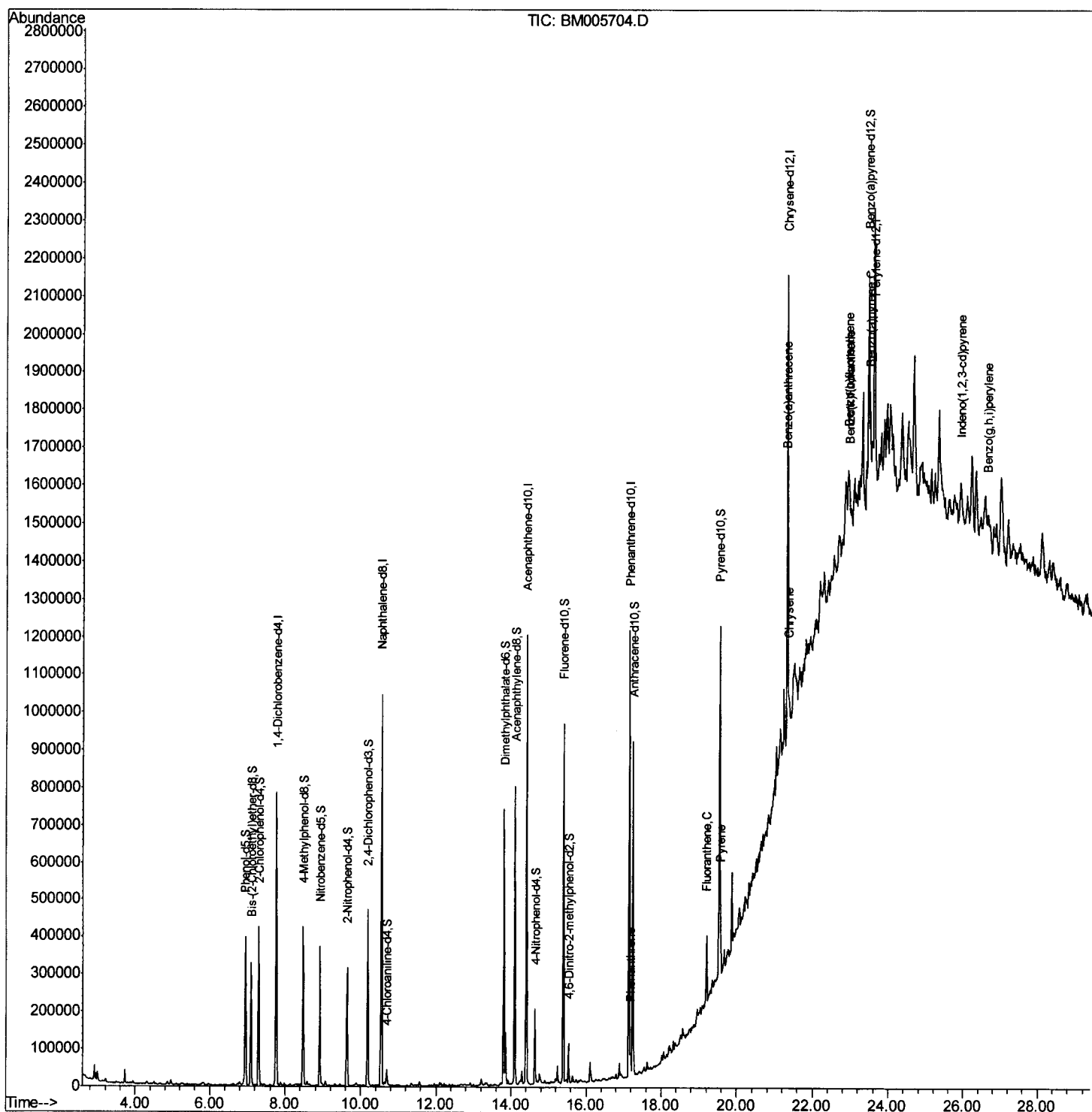
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060716\
 Data File : BM005704.D
 Acq On : 07 Jun 2016 22:15
 Operator : UM/SJ
 Sample : H3412-19
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9XT8

Manual Integrations
 APPROVED

sohil
 6/8/2016 7:16:24 PM

Quant Time: Jun 08 02:00:58 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM060216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 08 01:24:32 2016
 Response via : Initial Calibration



Quantitation Report (Qedit)

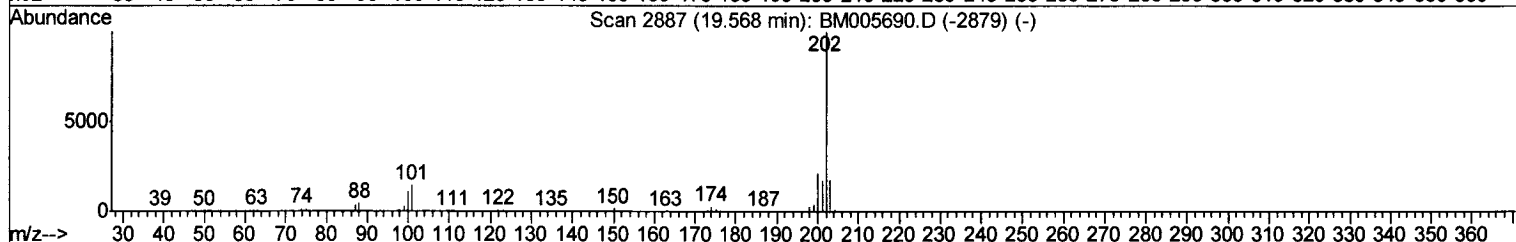
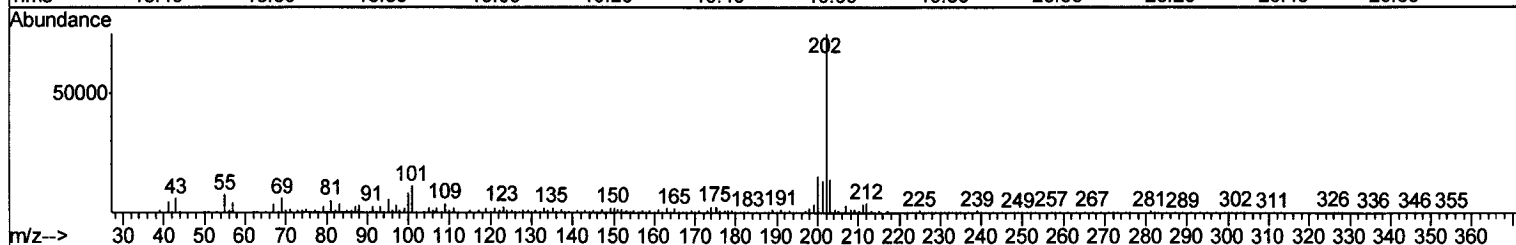
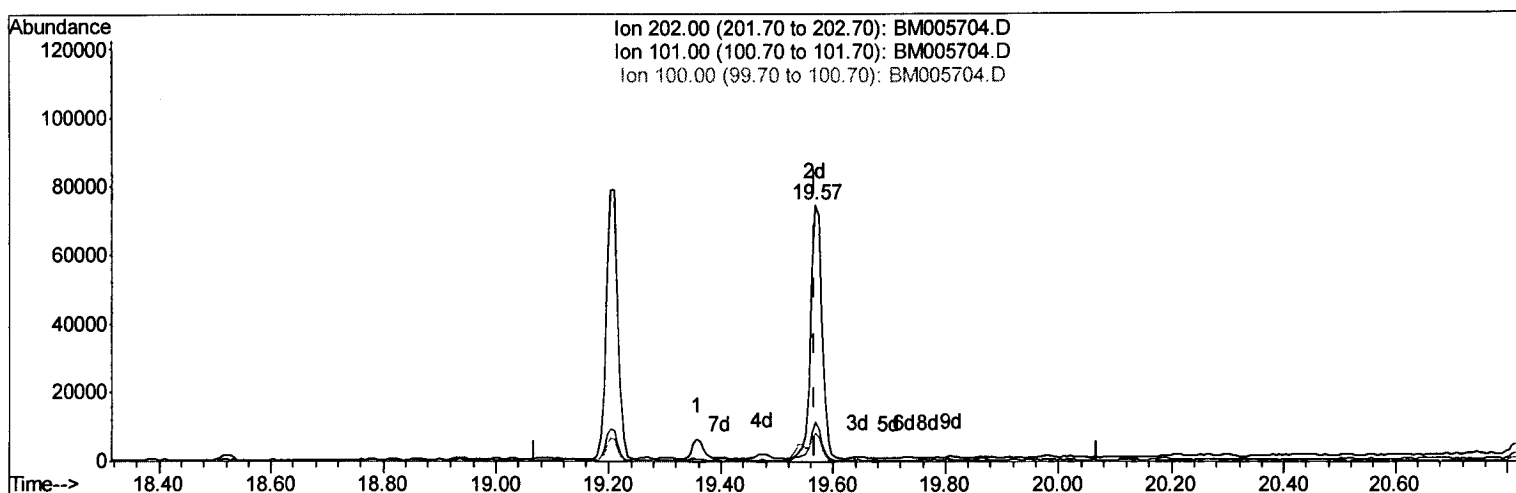
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TIC: BM005704.D

(31) Pyrene

19.568min (+0.000) 3.01ng/ul m

response 112132

Ion	Exp%	Act%
202.00	100	100
101.00	12.30	15.20#
100.00	11.60	11.12
0.00	0.00	0.00

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Quantitation Report (Qedit)

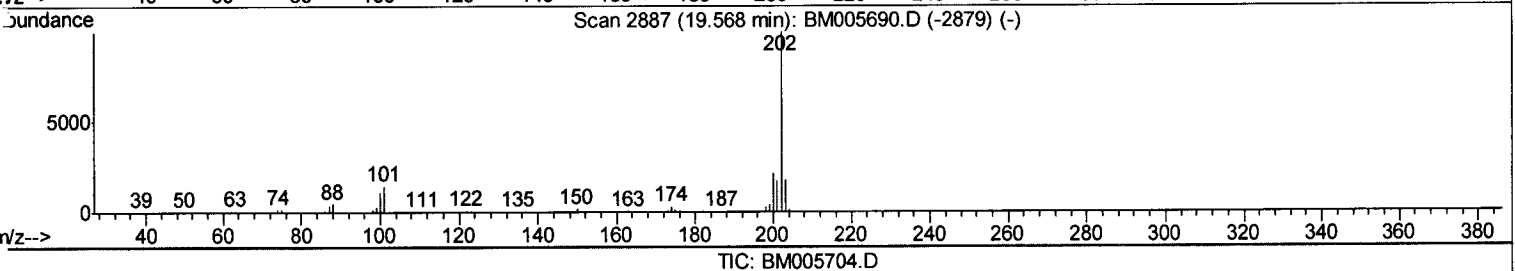
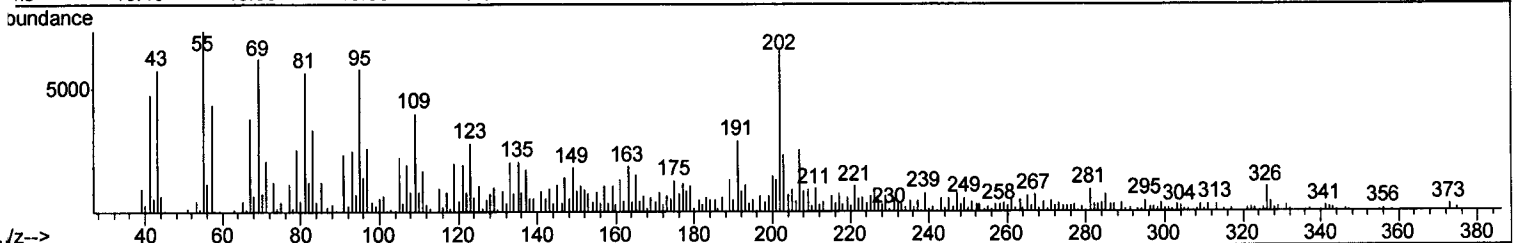
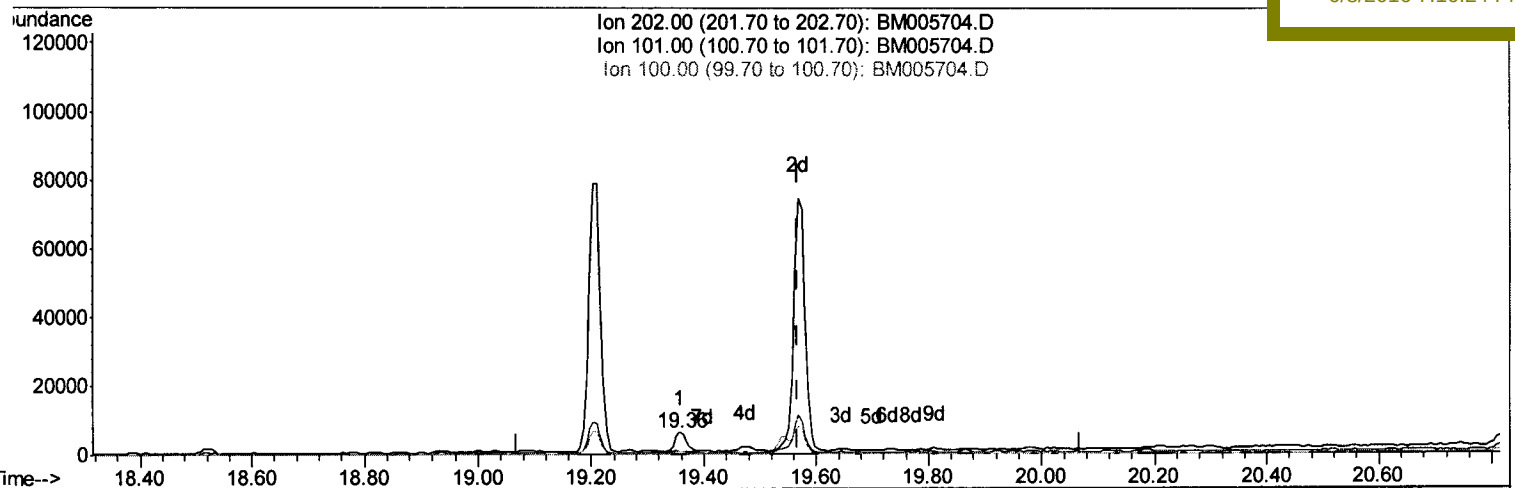
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(31) Pyrene

19.356min (-0.212) 0.22ng/ul

response 8157

Ion	Exp%	Act%
202.00	100	100
101.00	12.30	13.87
100.00	11.60	12.51
0.00	0.00	0.00

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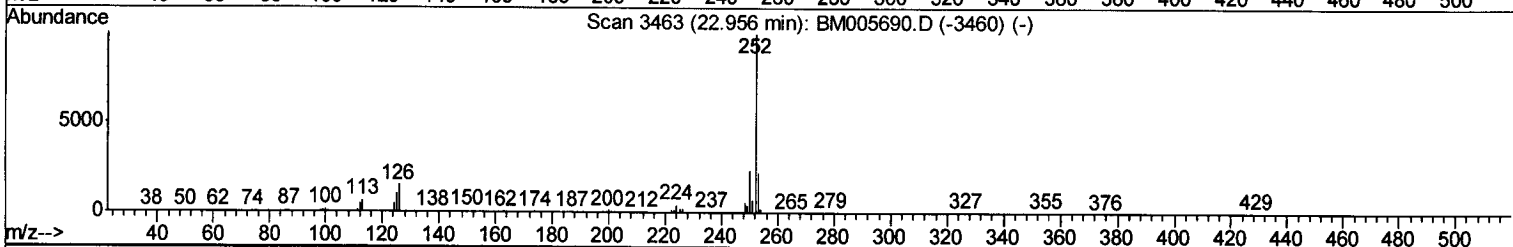
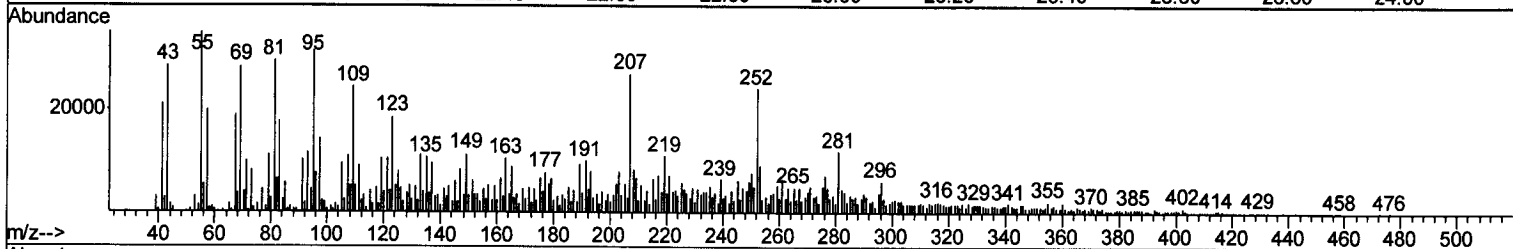
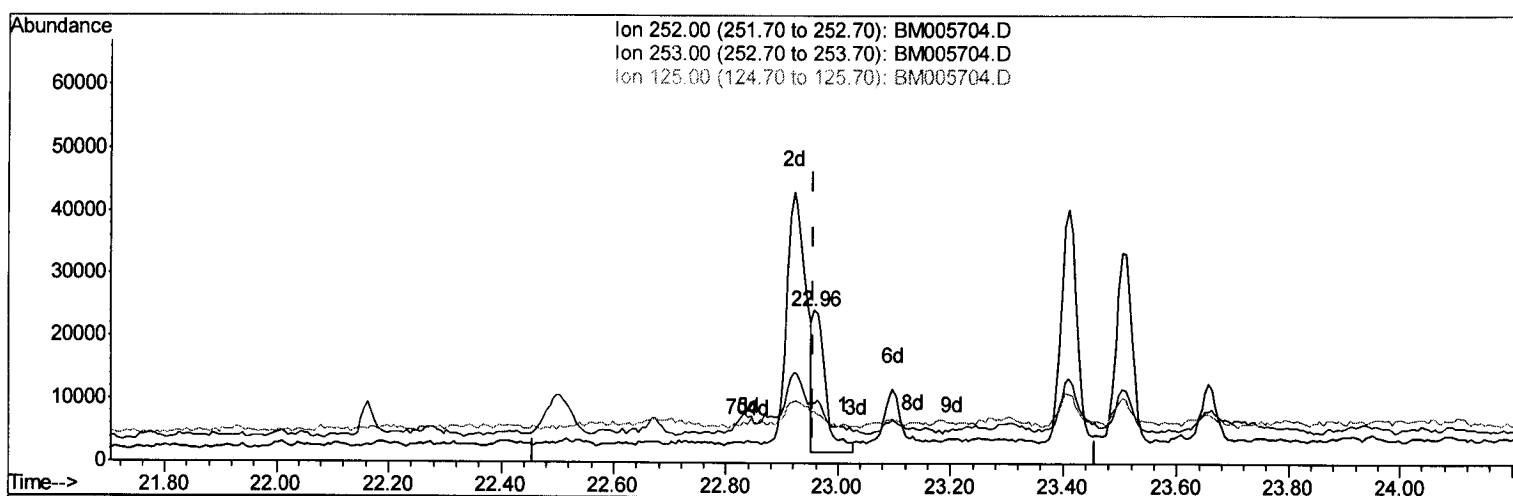
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TIC: BM005704.D

(36) Benzo(k)fluoranthene

22.956min (+0.000) 1.08ng/ul m

response 36243

Ion Exp% Act%

252.00 100 100

253.00 22.30 38.55#

125.00 9.60 33.92#

0.00 0.00 0.00

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Quantitation Report (Qedit)

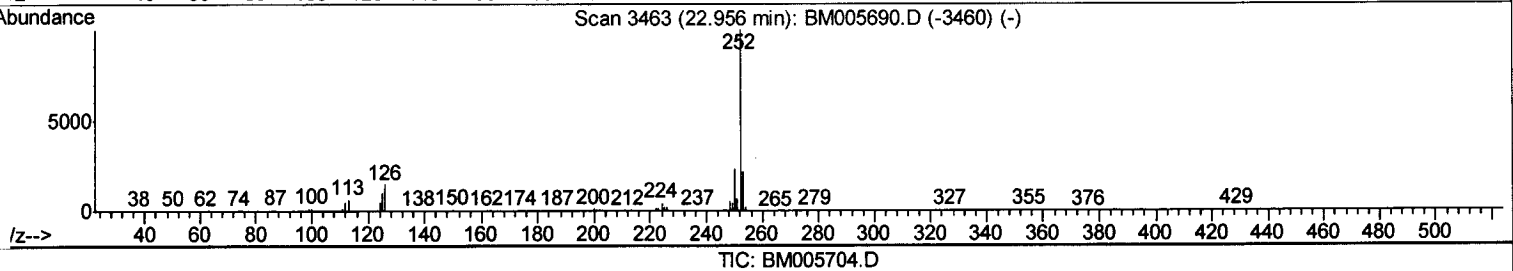
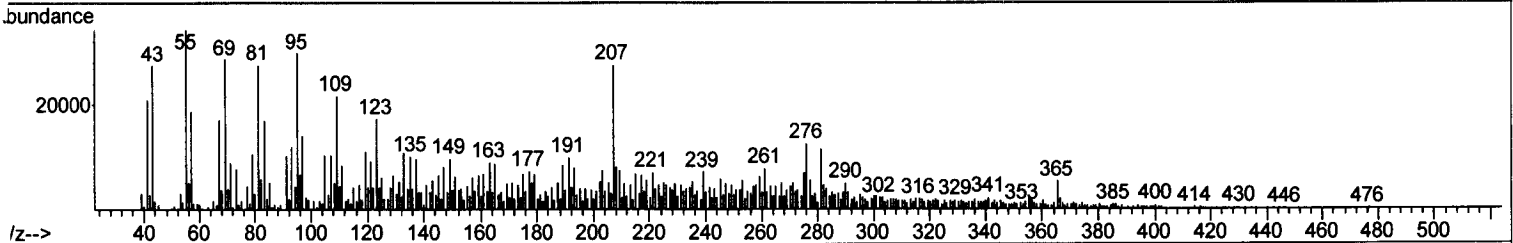
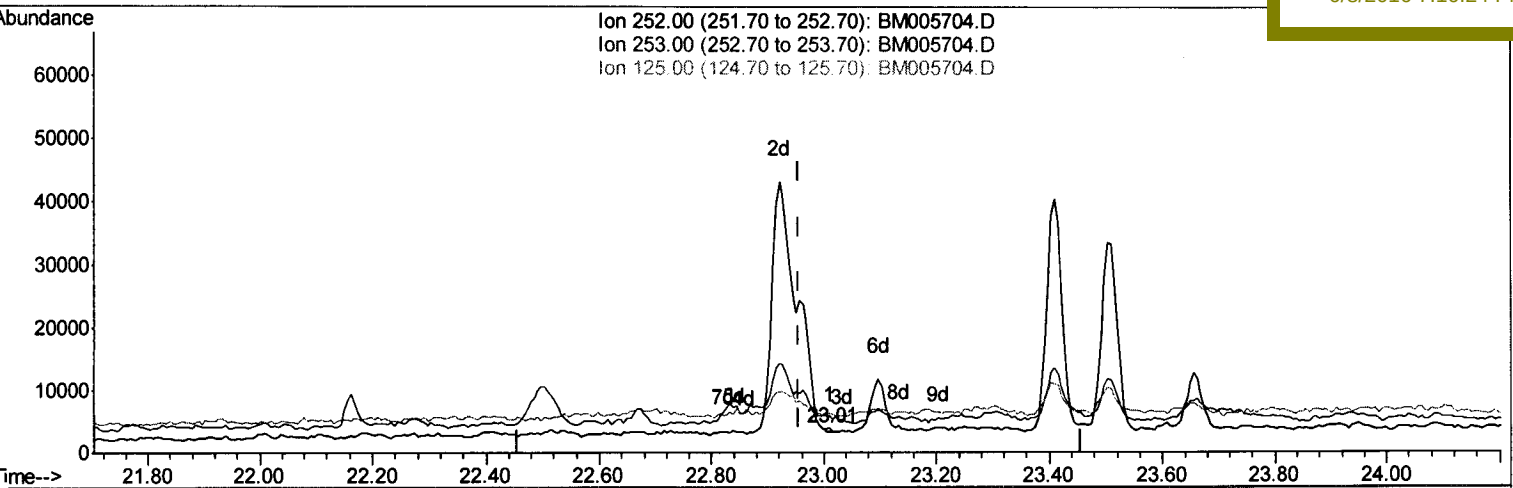
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(36) Benzo(k)fluoranthene

23.009min (+0.053) 0.01ng/ul

response 307

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	147.03#
125.00	9.60	159.31#
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	212695	20.00	ng/ul	0.00
7) Naphthalene-d8	10.55	136	879668	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.40	164	400325	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.15	188	696814	20.00	ng/ul	0.00
29) Chrysene-d12	21.33	240	590999	20.00	ng/ul	0.00
34) Perylene-d12	23.61	264	591863	20.00	ng/ul	0.01

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.21	96	4749	0.94	ng/uL	0.00
3) Phenol-d5	6.95	99	239267	12.80	ng/ul	0.00
4) Bis-(2-Chloroethyl) ether-d	7.09	67	142382	12.51	ng/ul	0.00
5) 2-Chlorophenol-d4	7.29	132	203708	13.20	ng/ul	0.00
6) 4-Methylphenol-d8	8.47	113	160870	10.88	ng/ul	0.00
8) Nitrobenzene-d5	8.92	128	95563	14.15	ng/ul	0.00
9) 2-Nitrophenol-d4	9.64	143	103375	14.30	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.19	165	177484	13.82	ng/ul	0.00
12) 4-Chloroaniline-d4	10.69	131	23986	1.57	ng/ul	0.00
16) Dimethylphthalate-d6	13.81	166	470499	15.29	ng/ul	0.00
17) Acenaphthylene-d8	14.10	160	584321	14.60	ng/ul	0.00
20) 4-Nitrophenol-d4	14.64	143	53625	12.17	ng/ul	0.00
21) Fluorene-d10	15.40	176	371910	14.07	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.54	200	20099	6.13	ng/ul	0.00
26) Anthracene-d10	17.24	188	491812	14.86	ng/ul	0.00
30) Pyrene-d10	19.54	212	466155	15.69	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.46	264	393081	14.43	ng/ul	0.01

Target Compounds

						Qvalue
25) Phenanthrene	17.19	178	47132	1.23	ng/ul	97
28) Fluoranthene	19.21	202	112479	2.86	ng/ul	99
31) Pyrene	19.57	202	112132m	3.01	ng/ul	
32) Benzo(a)anthracene	21.32	228	59230	1.71	ng/ul#	85
33) Chrysene	21.37	228	63003	1.92	ng/ul#	89
35) Benzo(b)fluoranthene	22.92	252	94003	2.63	ng/ul#	74
36) Benzo(k)fluoranthene	22.96	252	36243m	1.08	ng/ul	
38) Benzo(a)pyrene	23.50	252	58452	1.70	ng/ul#	63
39) Indeno(1,2,3-cd)pyrene	25.92	276	50827	1.25	ng/ul#	67
41) Benzo(g,h,i)perylene	26.64	276	62127	1.82	ng/ul#	74

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

UM
 06/08/16