

Quantitation Report (QT Reviewed)

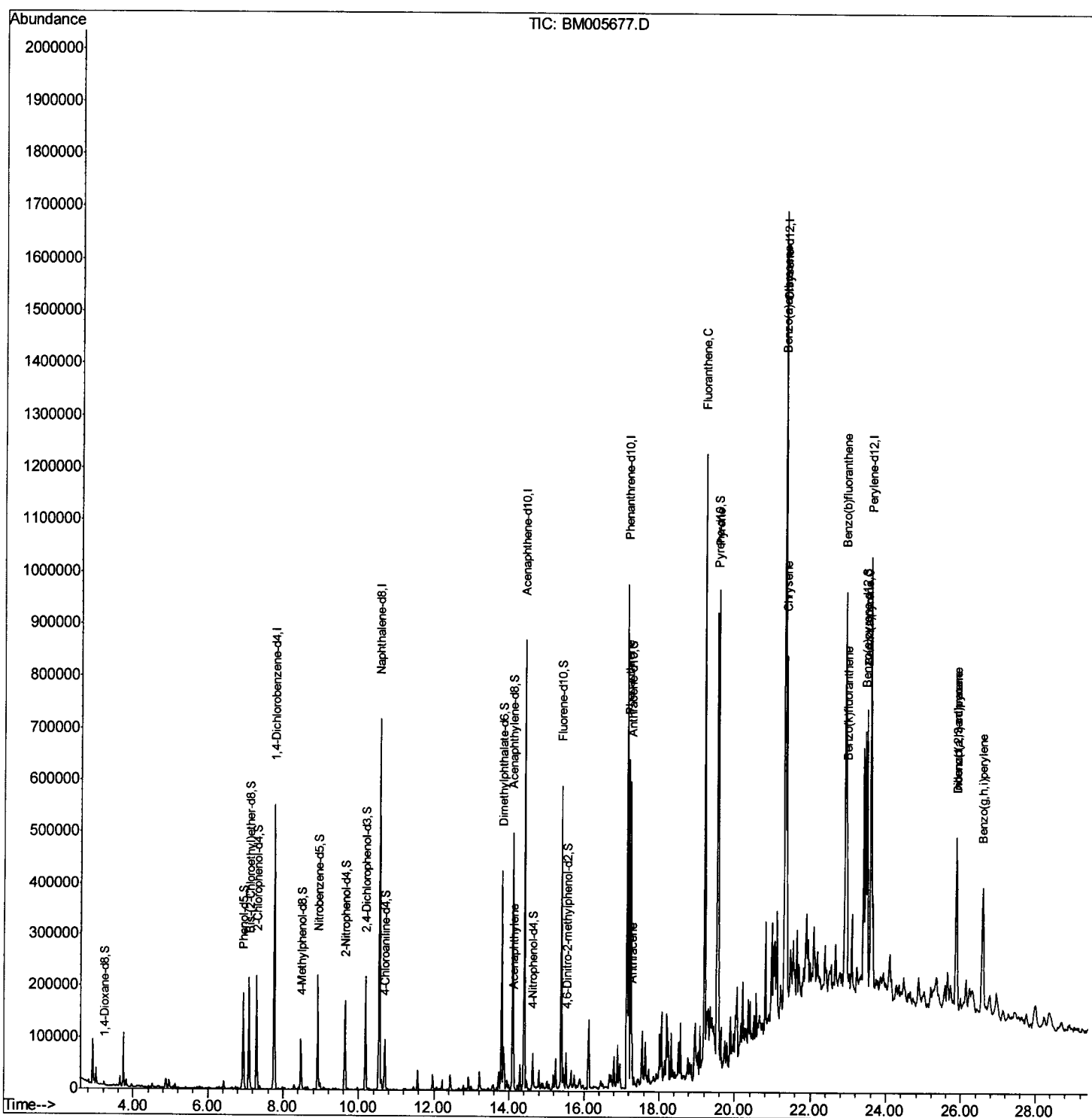
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060616\
 Data File : BM005677.D
 Acq On : 06 Jun 2016 19:36
 Operator : UM/SJ
 Sample : H3412-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9XS5

Manual Integrations
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Quant Time: Jun 07 04:57:39 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM060216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 07 04:25:20 2016
 Response via : Initial Calibration



Quantitation Report (Qedit)

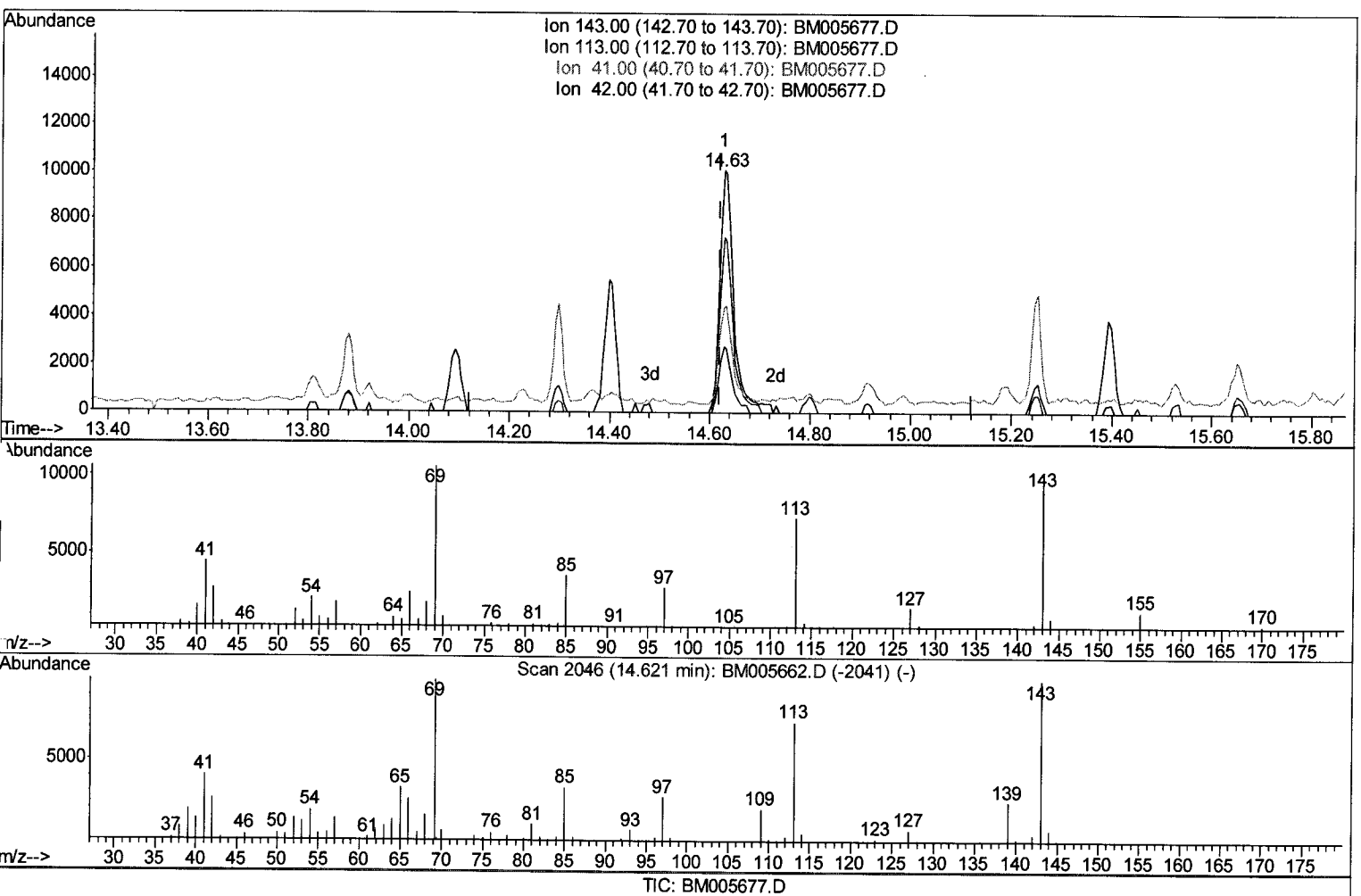
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(20) 4-Nitrophenol-d4 (S)

14.627min (+0.006) 6.39ng/ul m

response 19960

Ion	Exp%	Act%
143.00	100	100
113.00	102.70	72.35#
41.00	49.70	43.85
42.00	33.70	27.13

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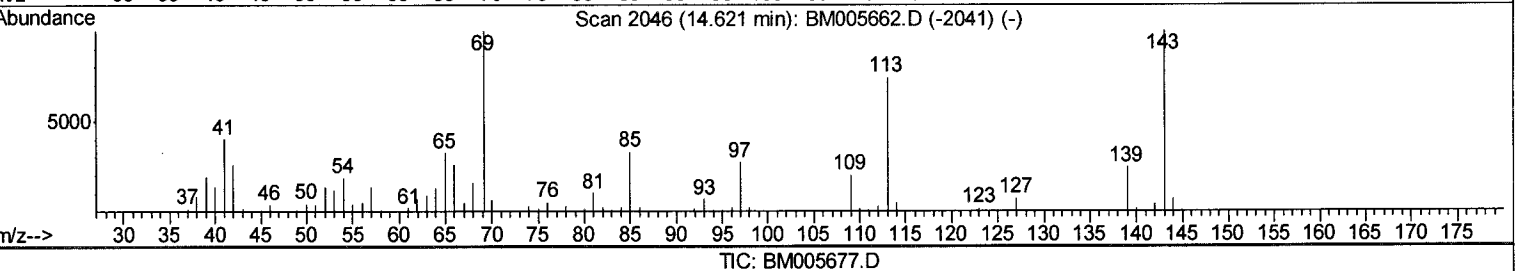
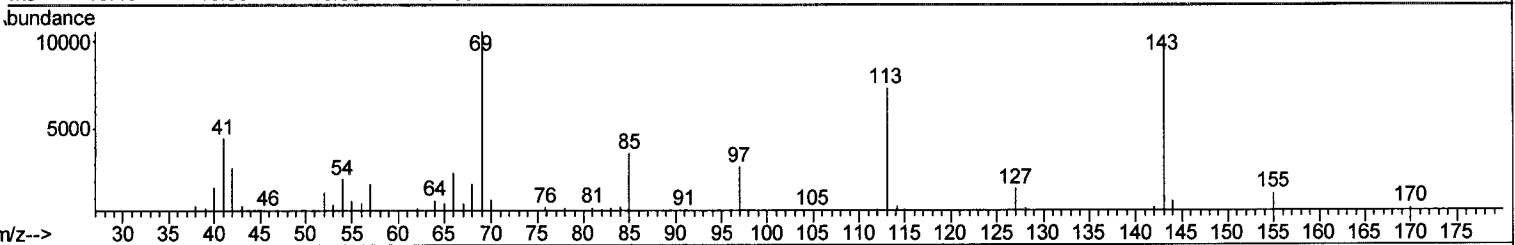
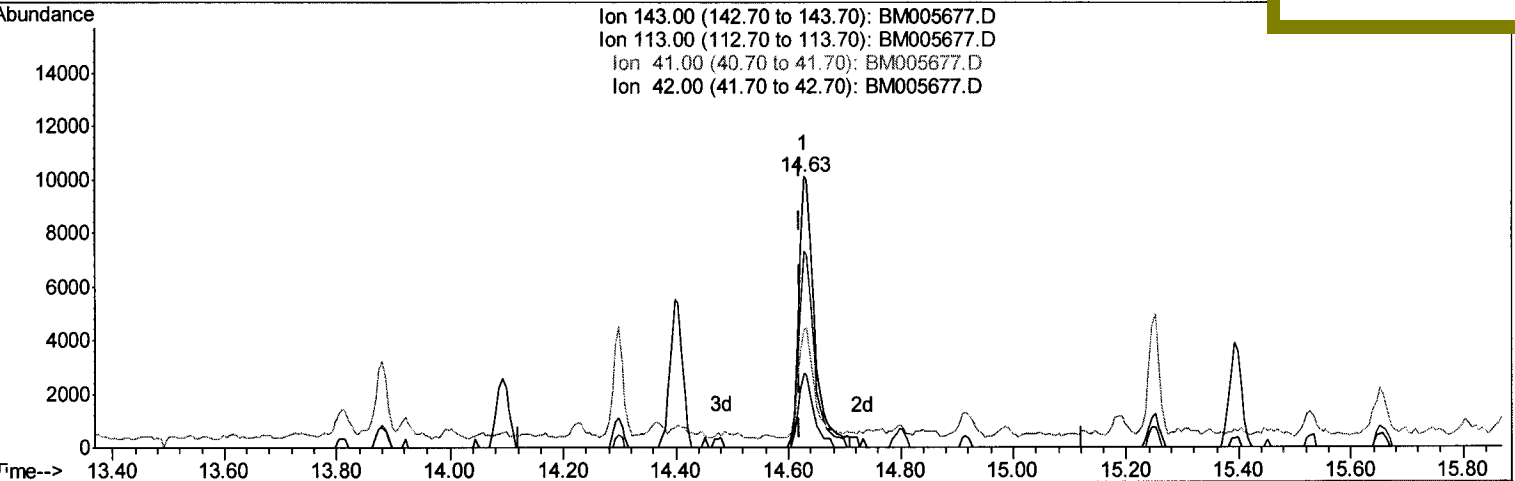
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(20) 4-Nitrophenol-d4 (S)

14.627min (+0.006) 6.31ng/ul

response 19703

Ion	Exp%	Act%
143.00	100	100
113.00	102.70	72.35#
41.00	49.70	43.85
42.00	33.70	27.13

Quantitation Report (Qedit)

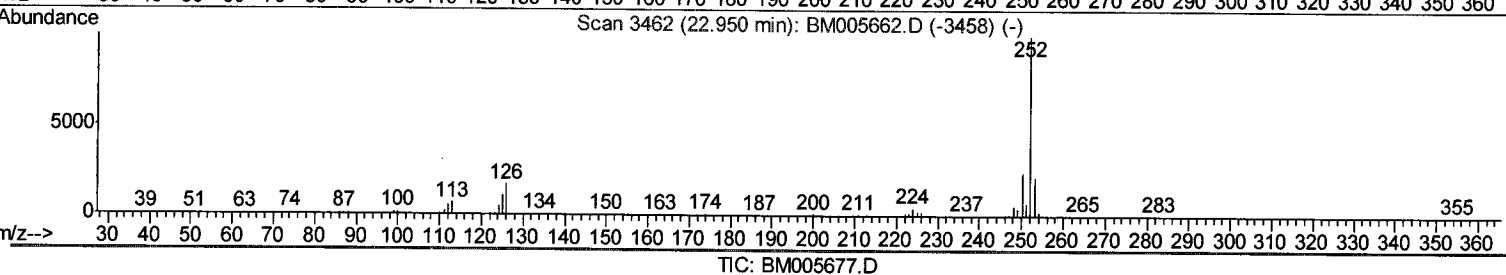
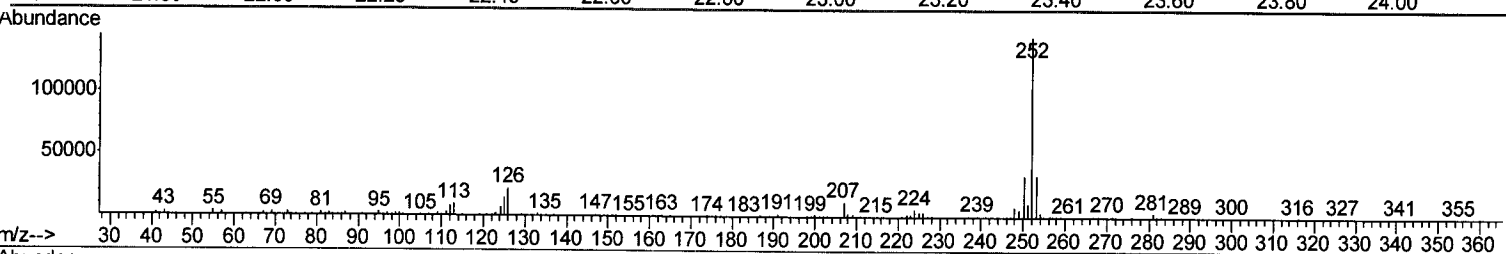
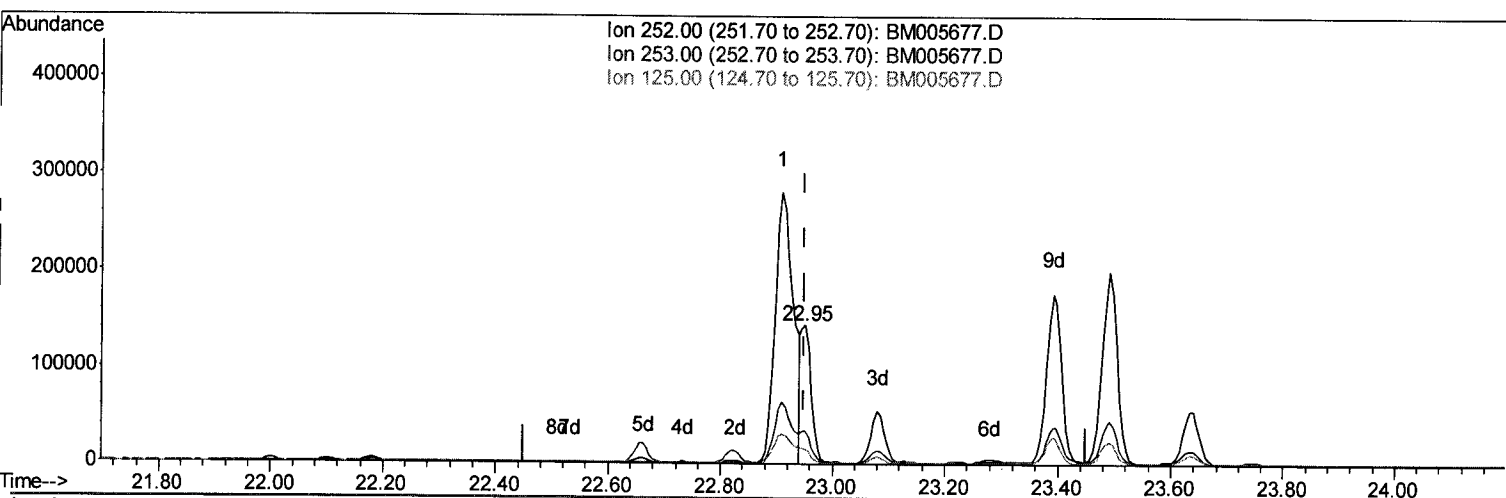
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TIC: BM005677.D

(36) Benzo(k)fluoranthene

22.950min (+0.000) 5.81ng/ul m

response 192539

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Ion	Exp%	Act%
252.00	100	100
253.00	22.30	23.59
125.00	9.60	10.81
0.00	0.00	0.00

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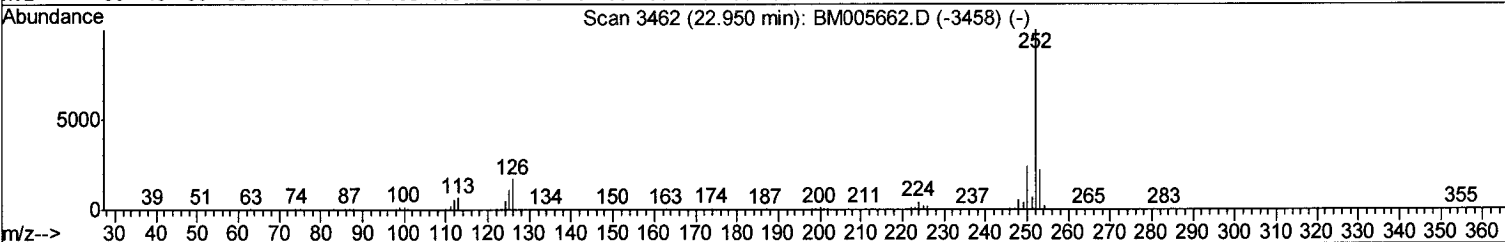
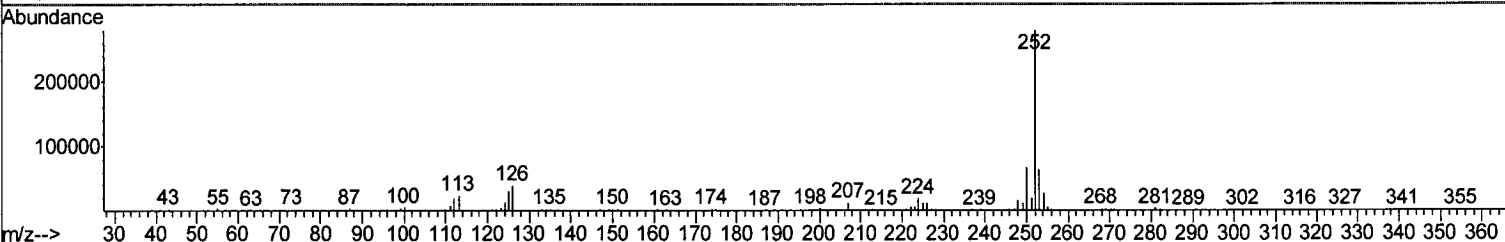
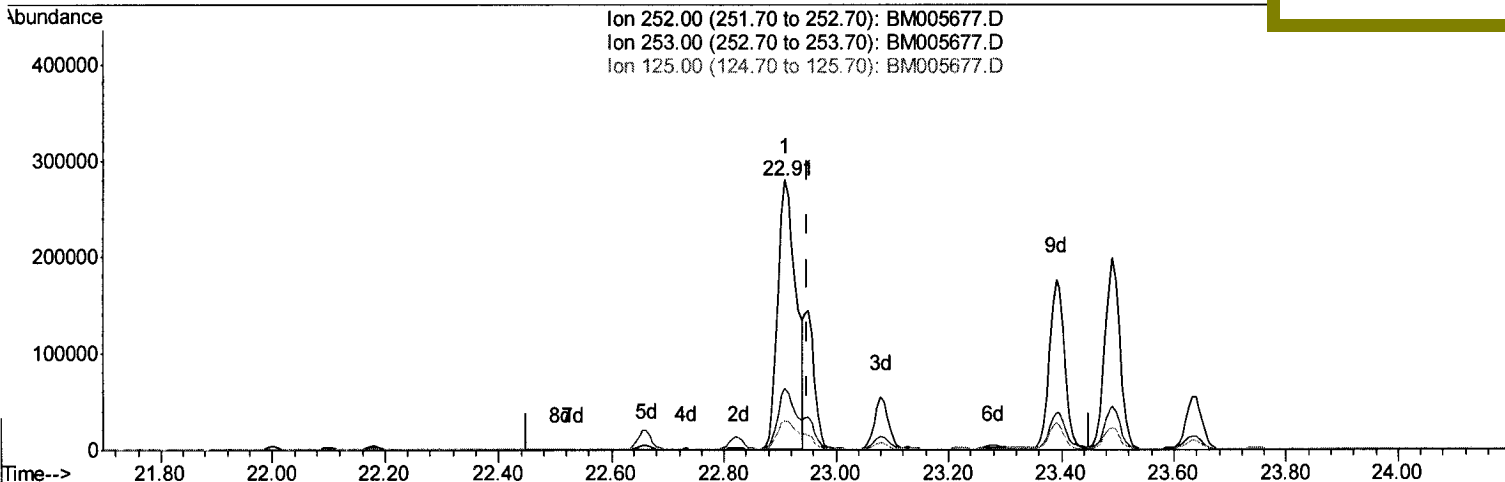
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TIC: BM005677.D

(36) Benzo(k)fluoranthene

22.909min (-0.041) 18.85ng/ul

response 624889

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	22.88
125.00	9.60	11.08
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	147920	20.00	ng/ul	0.00
7) Naphthalene-d8	10.55	136	599095	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.40	164	283843	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.14	188	556433	20.00	ng/ul	0.00
29) Chrysene-d12	21.33	240	556515	20.00	ng/ul	0.00
34) Perylene-d12	23.59	264	586057	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.22	96	3665	1.04	ng/uL	0.00
3) Phenol-d5	6.94	99	111076	8.54	ng/ul	0.00
4) Bis-(2-Chloroethyl) ether-d	7.09	67	93496	11.81	ng/ul	0.00
5) 2-Chlorophenol-d4	7.29	132	105267	9.81	ng/ul	0.00
6) 4-Methylphenol-d8	8.47	113	38650	3.76	ng/ul	0.00
8) Nitrobenzene-d5	8.92	128	53426	11.61	ng/ul	0.00
9) 2-Nitrophenol-d4	9.64	143	55921	11.36	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.18	165	81414	9.31	ng/ul	0.00
12) 4-Chloroaniline-d4	10.69	131	56554	5.44	ng/ul	0.00
16) Dimethylphthalate-d6	13.81	166	269709	12.36	ng/ul	0.00
17) Acenaphthylene-d8	14.09	160	347182	12.23	ng/ul	0.00
20) 4-Nitrophenol-d4	14.63	143	19960m	6.39	ng/ul	0.00
21) Fluorene-d10	15.39	176	231437	12.35	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	13930	5.32	ng/ul	0.00
26) Anthracene-d10	17.24	188	319542	12.09	ng/ul	0.00
30) Pyrene-d10	19.54	212	340145	12.16	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.44	264	320867	11.90	ng/ul	0.00

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Target Compounds

						Qvalue
18) Acenaphthylene	14.12	152	64382	2.19	ng/ul	97
25) Phenanthrene	17.19	178	366136	11.92	ng/ul	98
27) Anthracene	17.28	178	73116	2.34	ng/ul	99
28) Fluoranthene	19.20	202	647190	20.62	ng/ul	98
31) Pyrene	19.57	202	545285	15.53	ng/ul	97
32) Benzo(a)anthracene	21.31	228	350410	10.74	ng/ul	97
33) Chrysene	21.36	228	370007	12.00	ng/ul	96
35) Benzo(b)fluoranthene	22.91	252	624889	17.68	ng/ul	98
36) Benzo(k)fluoranthene	22.95	252	192539m	5.81	ng/ul	98
38) Benzo(a)pyrene	23.49	252	367312	10.80	ng/ul	98
39) Indeno(1,2,3-cd)pyrene	25.87	276	317114	7.88	ng/ul	96
40) Dibenzo(a,h)anthracene	25.88	278	88146	2.61	ng/ul#	92
41) Benzo(g,h,i)perylene	26.58	276	304282	9.01	ng/ul#	94

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