

Quantitation Report (QT Reviewed)

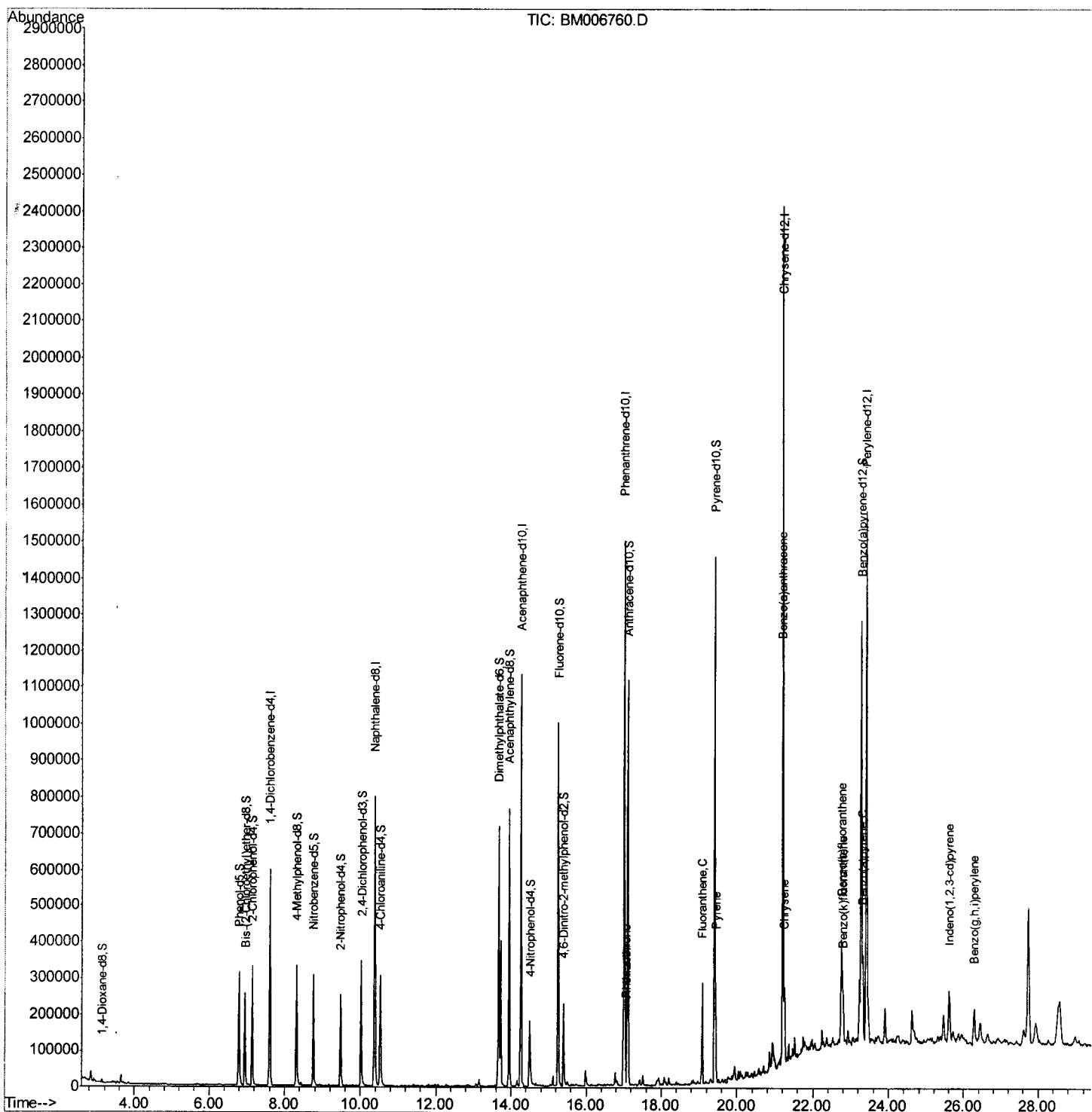
Data Path : Z:\HPCHEM1\BNA_M\Data\BM073016\
 Data File : BM006760.D
 Acq On : 30 Jul 2016 06:21
 Operator : UM/SJ
 Sample : H4188-03
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9ZL1

Manual Integrations
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Quant Time: Aug 01 04:53:15 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 30 01:37:26 2016
 Response via : Initial Calibration



Quantitation Report (Qedit)

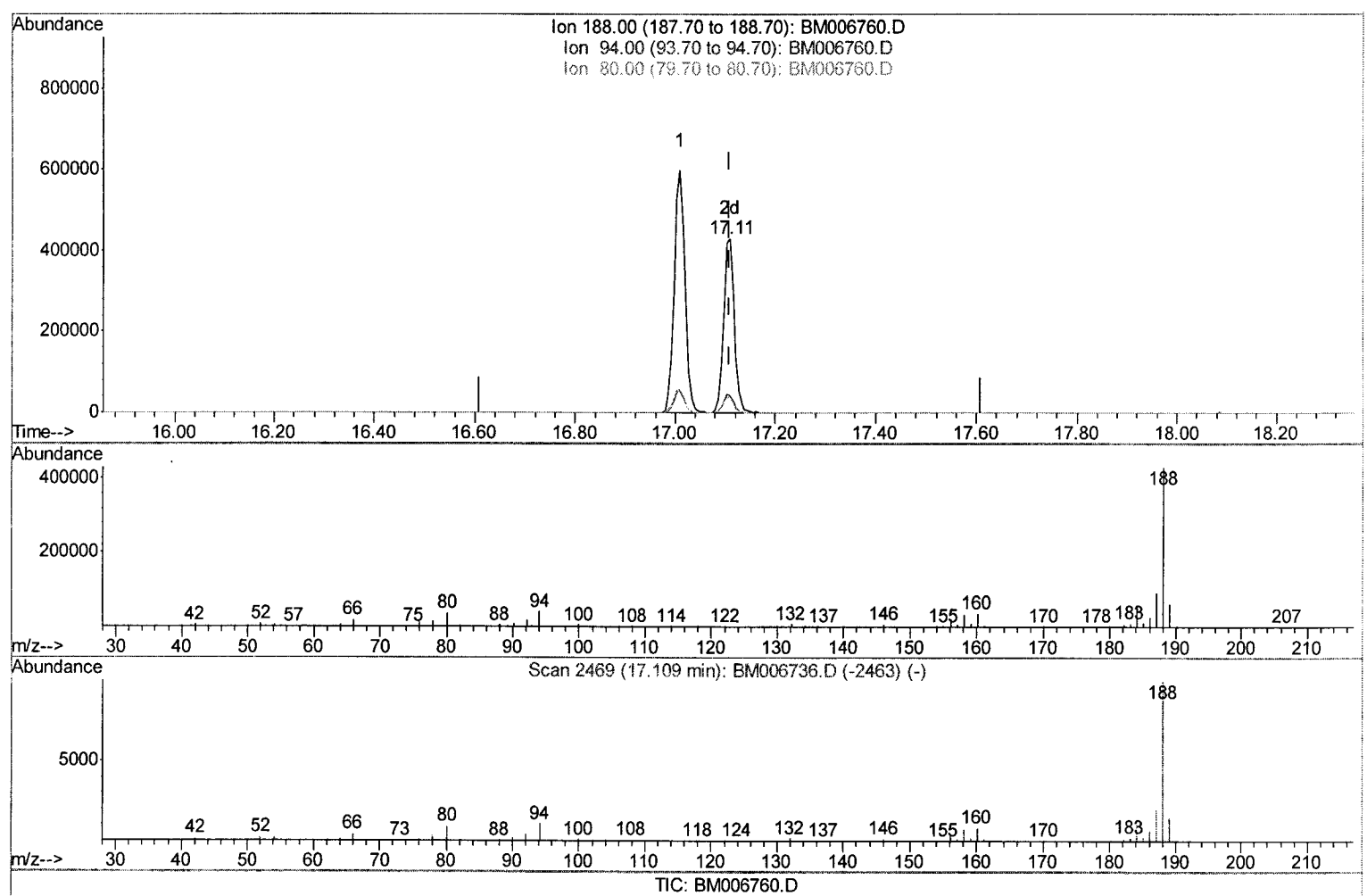
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(26) Anthracene-d10 (S)

17.110min (+0.000) 15.45ng/ul m

response 639363

Ion	Exp%	Act%
188.00	100	100
94.00	10.70	9.56
80.00	10.80	8.61#
0.00	0.00	0.00

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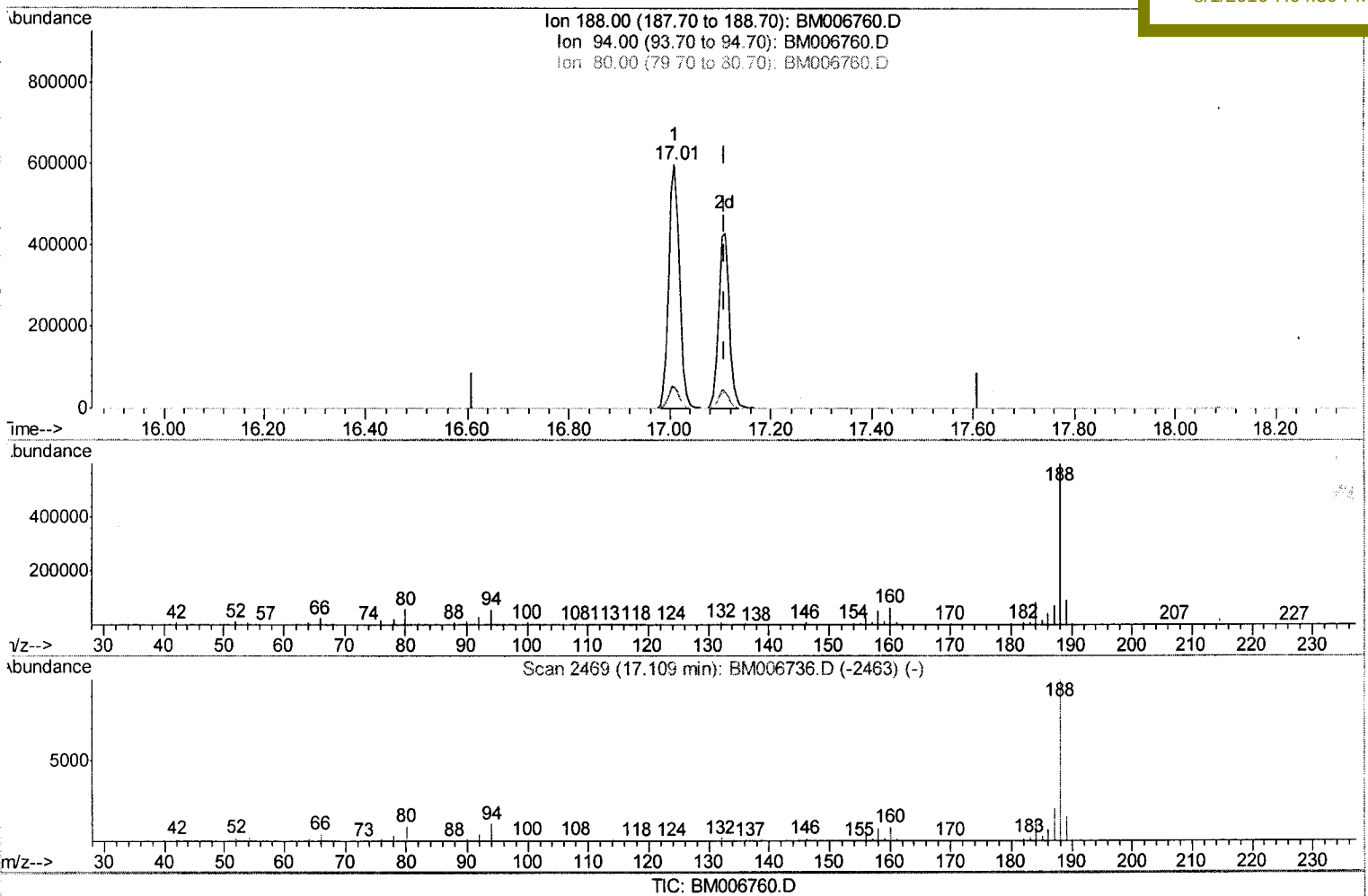
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(26) Anthracene-d10 (S)

17.010min (-0.100) 21.06ng/ul

response 871549

Ion	Exp%	Act%
188.00	100	100
94.00	10.70	8.97
80.00	10.80	9.43
0.00	0.00	0.00

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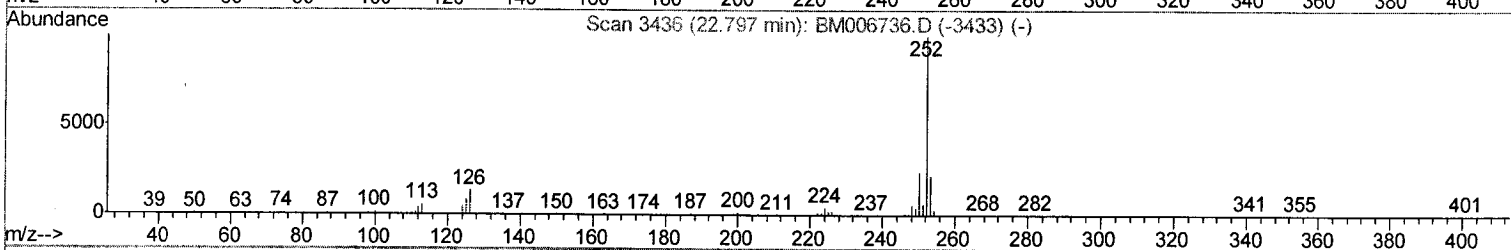
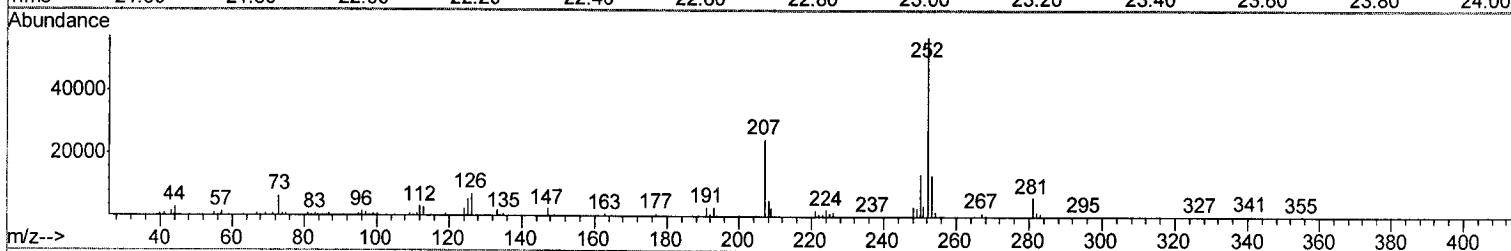
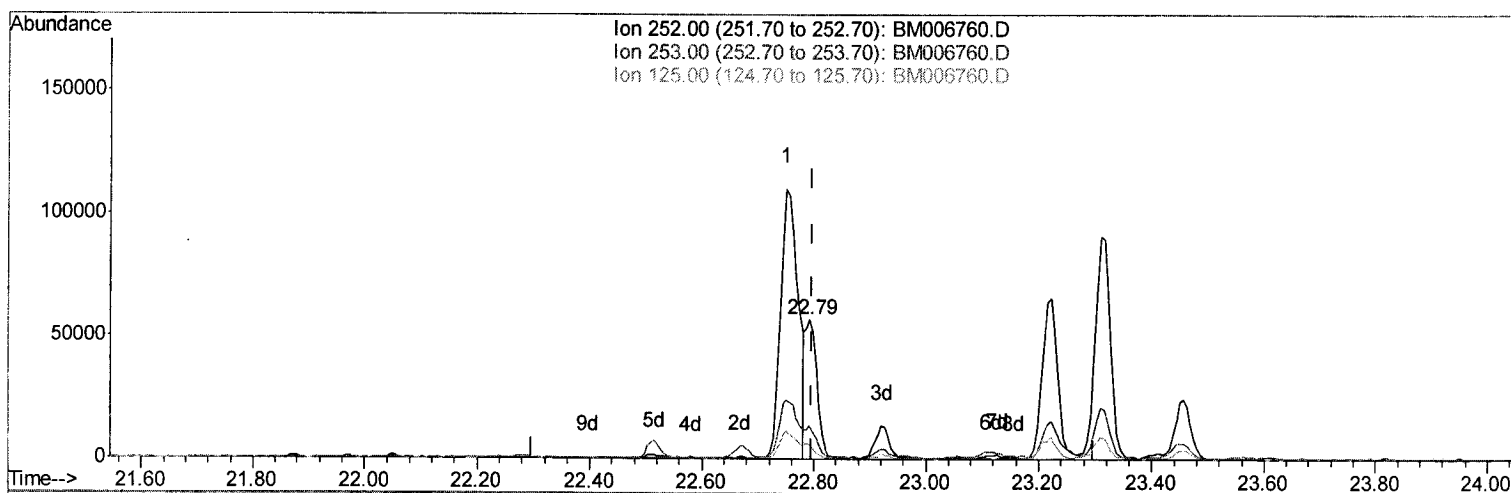
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Quant Time: Aug 01 04:51:54 2016
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TIC: BM006760.D

(36) Benzo(k)fluoranthene

22.791min (-0.006) 1.31ng/ul m *U.M*
08/02/16
 response 83372

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	23.77
125.00	9.60	10.22
0.00	0.00	0.00

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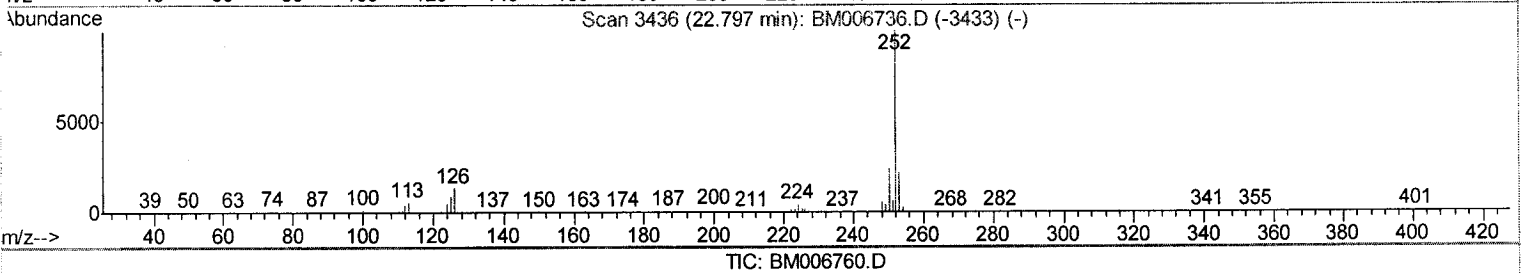
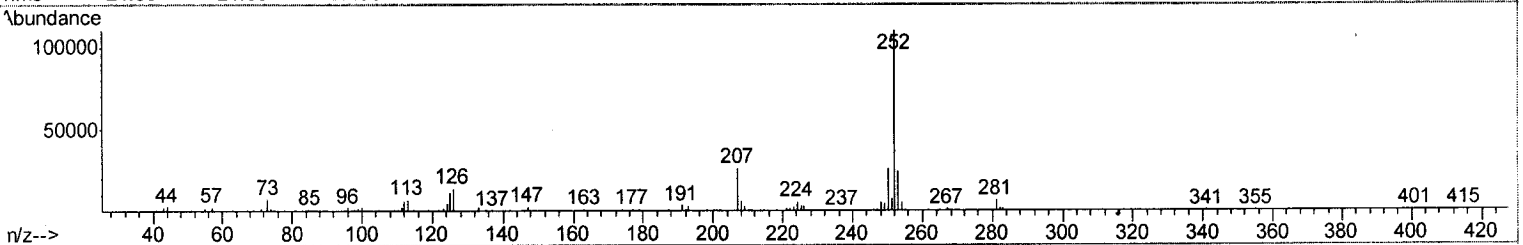
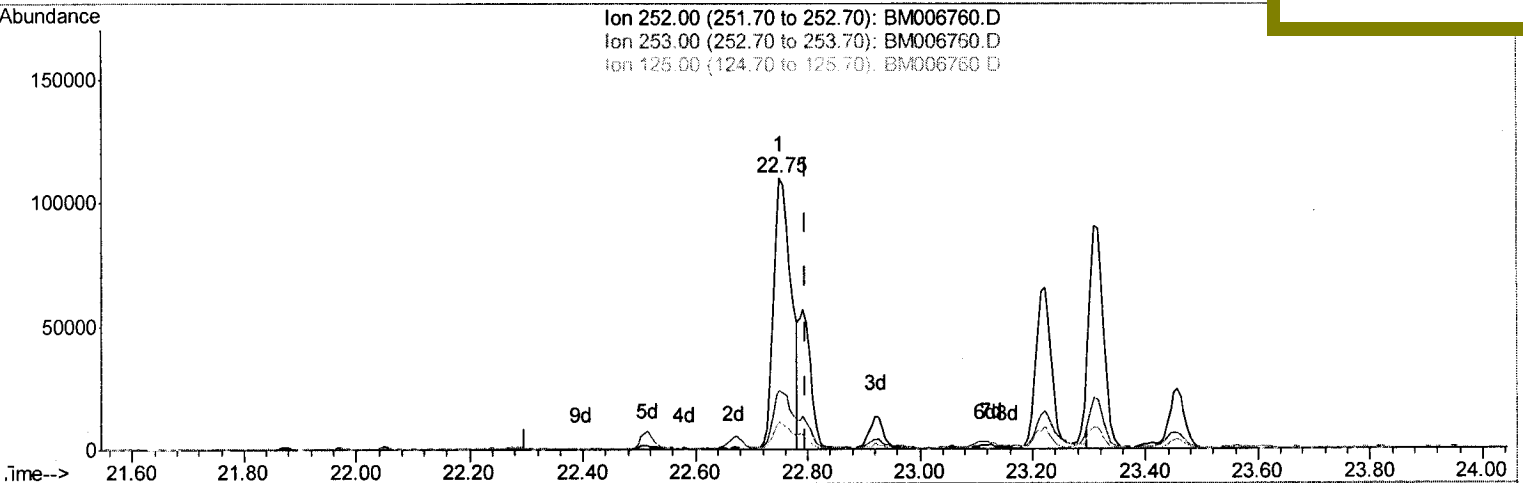
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(36) Benzo(k)fluoranthene
 22.750min (-0.047) 3.76ng/ul
 response 239152

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	21.89
125.00	9.60	10.47
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	7.60	152	154703	20.00	ng/ul	0.00
7) Naphthalene-d8	10.38	136	664113	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.26	164	380469	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.01	188	871549	20.00	ng/ul	0.00
29) Chrysene-d12	21.21	240	1053203	20.00	ng/ul	0.00
34) Perylene-d12	23.41	264	1098696	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.13	96	4889	1.28	ng/uL	0.00
3) Phenol-d5	6.79	99	173091	13.15	ng/ul	0.00
4) Bis-(2-Chloroethyl) ether-d	6.95	67	120340	14.76	ng/ul	0.00
5) 2-Chlorophenol-d4	7.14	132	141294	13.44	ng/ul	0.00
6) 4-Methylphenol-d8	8.32	113	127540	12.49	ng/ul	0.00
8) Nitrobenzene-d5	8.76	128	68966	13.72	ng/ul	0.00
9) 2-Nitrophenol-d4	9.47	143	77251	14.45	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.02	165	137513	13.61	ng/ul	0.00
12) 4-Chloroaniline-d4	10.53	131	176059	15.82	ng/ul	0.00
16) Dimethylphthalate-d6	13.67	166	463698	15.54	ng/ul	0.00
17) Acenaphthylene-d8	13.95	160	567568	14.91	ng/ul	0.00
20) 4-Nitrophenol-d4	14.49	143	53431	10.31	ng/ul	0.00
21) Fluorene-d10	15.26	176	414217	15.53	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.40	200	48599	10.74	ng/ul	0.00
26) Anthracene-d10	17.11	188	639363m	15.45	ng/ul	0.00
30) Pyrene-d10	19.41	212	751922	15.67	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.27	264	786372	15.56	ng/ul	-0.01

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Target Compounds

					Qvalue
25) Phenanthrene	17.05	178	72919	1.49 ng/ul	99
27) Anthracene	17.05	178	72919	1.44 ng/ul	97
28) Fluoranthene	19.07	202	162475	2.82 ng/ul	99
31) Pyrene	19.44	202	147927	2.32 ng/ul	98
32) Benzo(a)anthracene	21.19	228	114323	1.79 ng/ul	98
33) Chrysene	21.24	228	132930	2.26 ng/ul	97
35) Benzo(b)fluoranthene	22.75	252	239152	3.63 ng/ul	98
36) Benzo(k)fluoranthene	22.79	252	83372m	1.31 ng/ul	98
38) Benzo(a)pyrene	23.31	252	164128	2.56 ng/ul	98
39) Indeno(1,2,3-cd)pyrene	25.60	276	132884	1.79 ng/ul	99
41) Benzo(g,h,i)perylene	26.28	276	116511	1.83 ng/ul	95

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