

Quantitation Report (QT Reviewed)

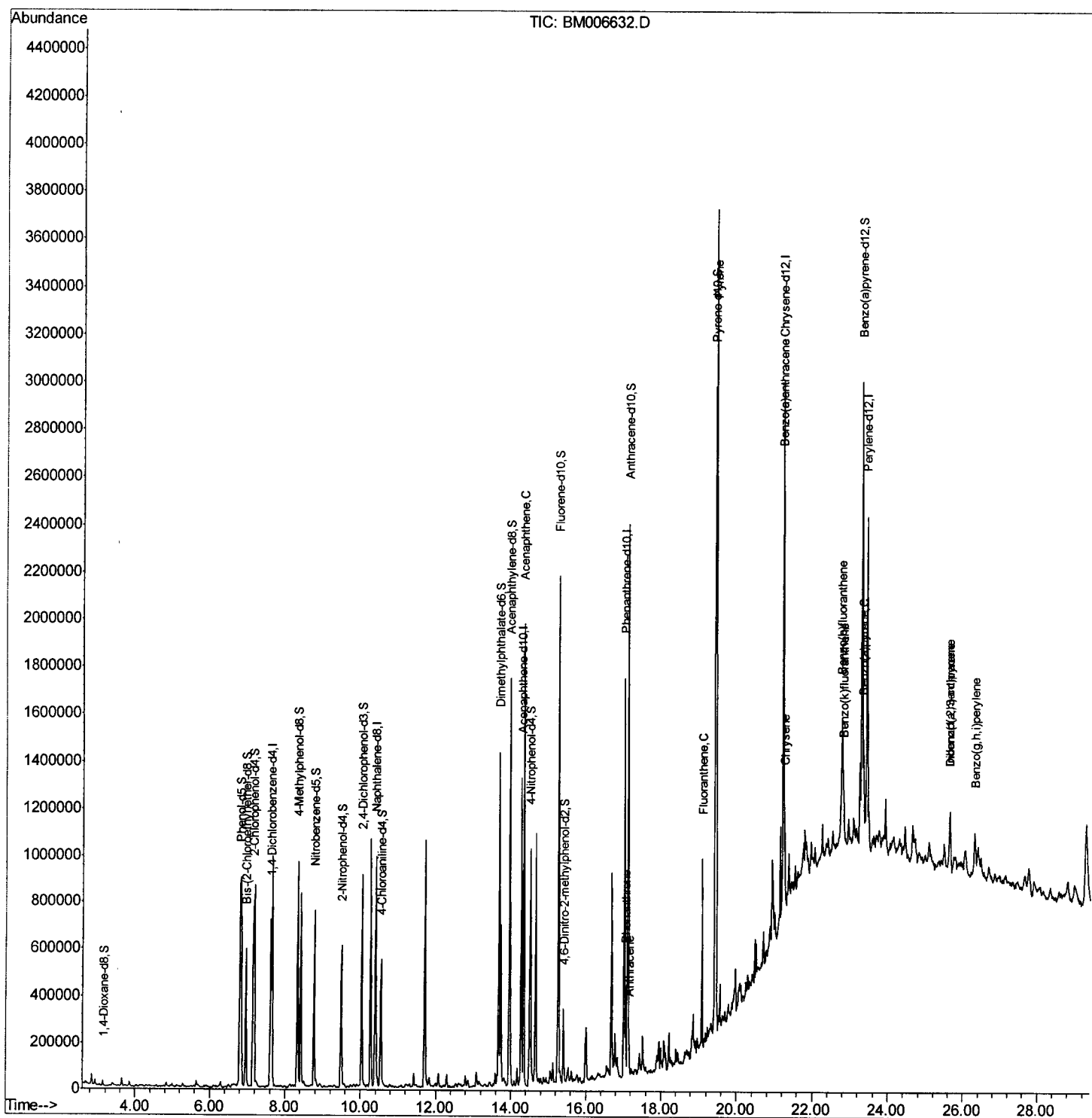
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM072216\
Data File : BM006632.D
Acq On : 22 Jul 2016 19:52
Operator : UM/SJ
Sample : H4076-14MSD
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
D9YP7MSD

Manual Integration

Quant Time: Jul 23 01:02:19 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jul 23 00:16:38 2016
Response via : Initial Calibration

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Quantitation Report (Qedit)

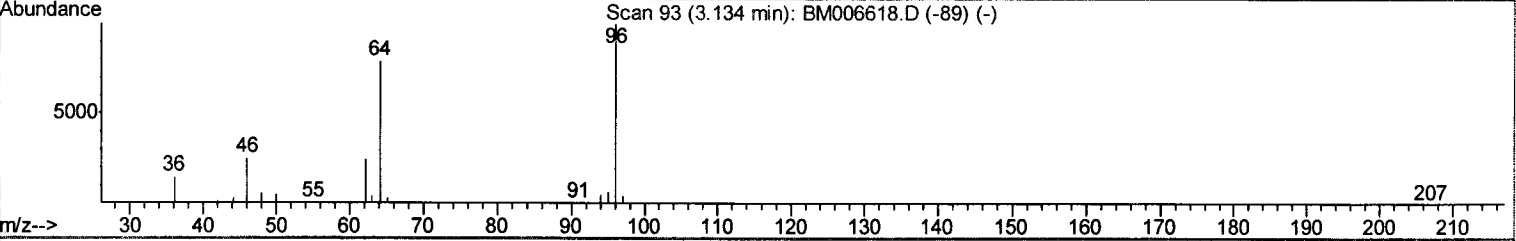
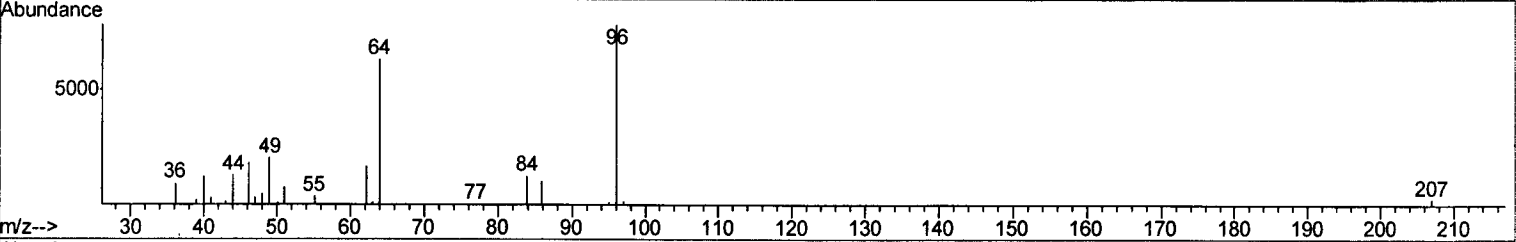
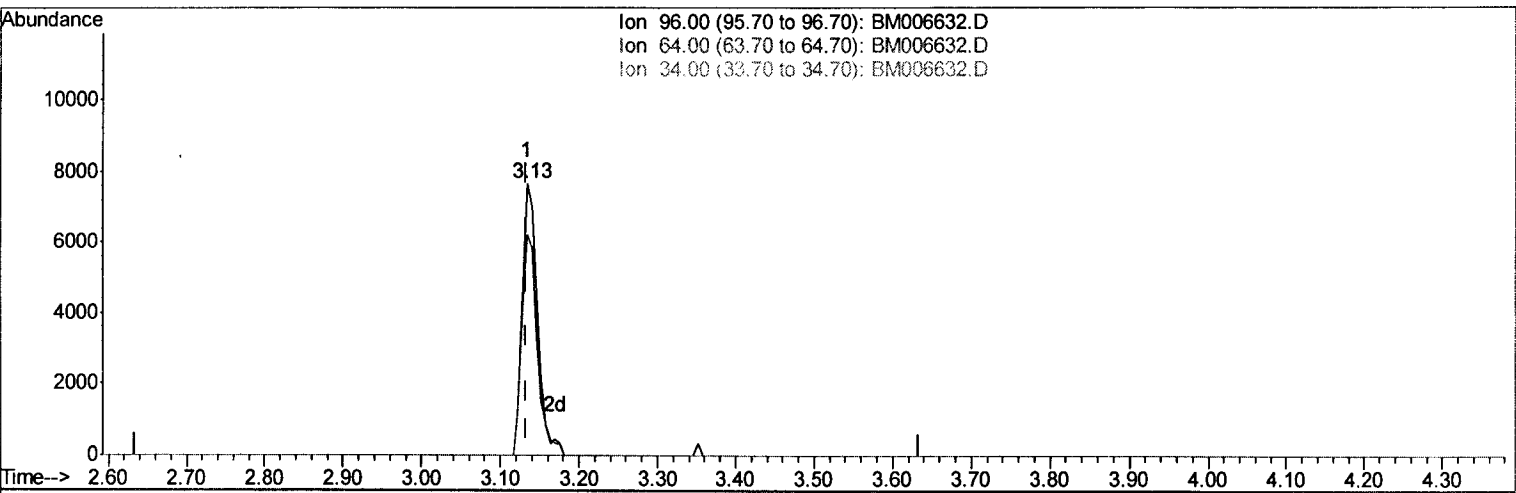
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TIC: BM006632.D

(2) 1,4-Dioxane-d8 (S)

3.134min (+0.000) 2.26ng/uL m

response 10211

Ion	Exp%	Act%
96.00	100	100
64.00	66.20	81.46#
34.00	0.00	0.00
0.00	0.00	0.00

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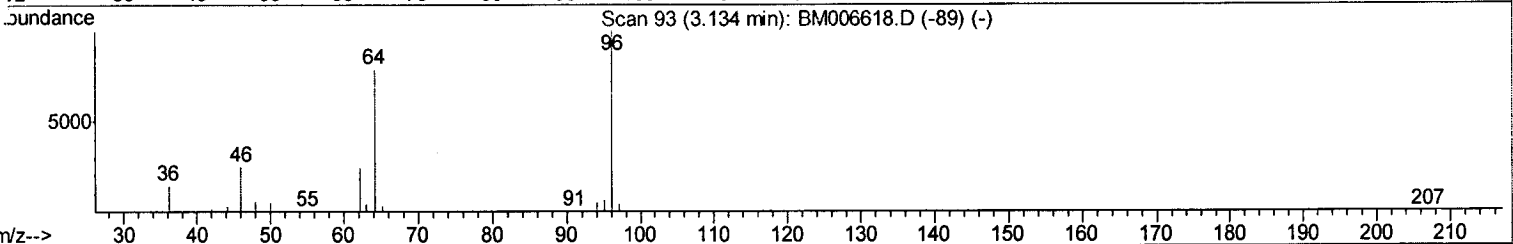
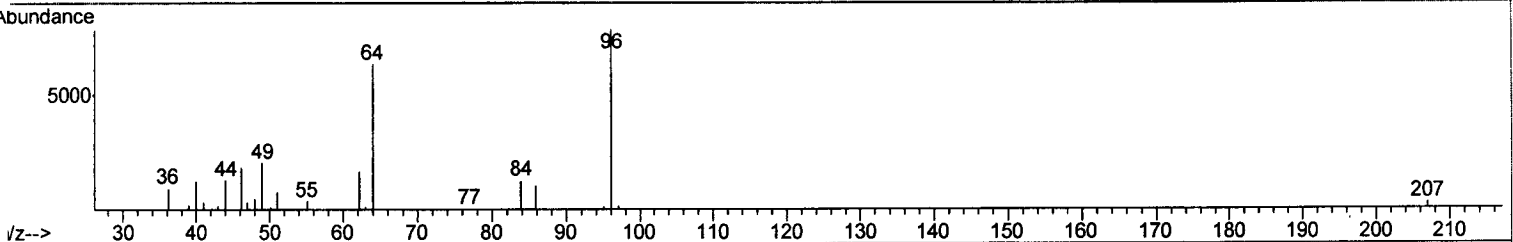
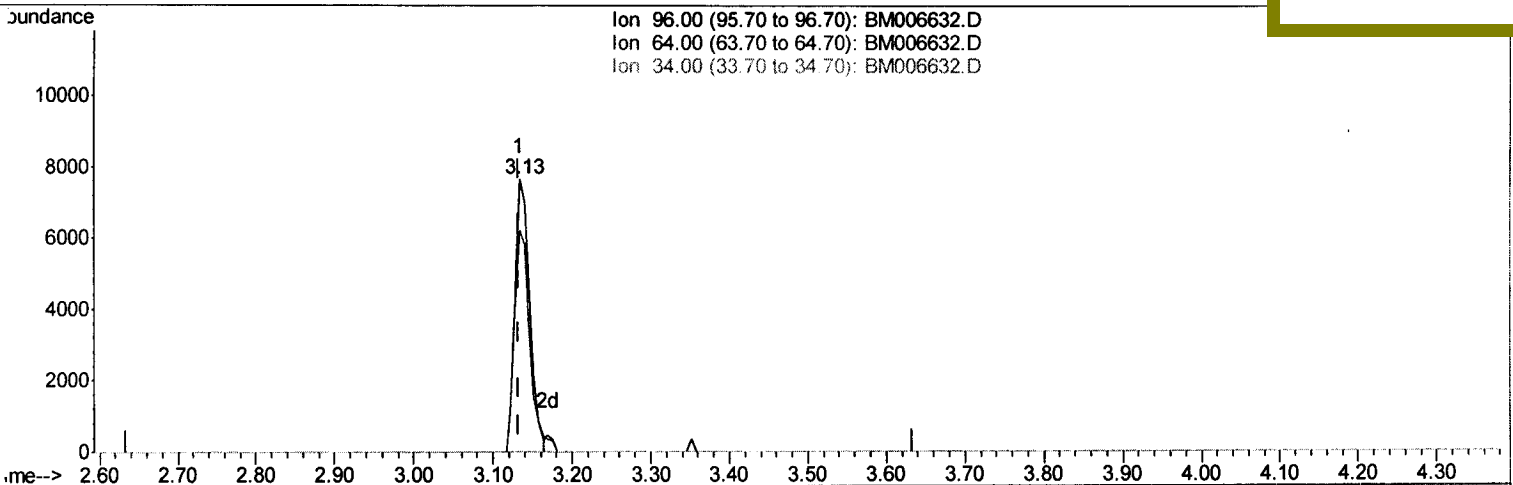
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TIC: BM006632.D

(2) 1,4-Dioxane-d8 (S)

3.134min (+0.000) 2.20ng/uL

response 9926

Ion	Exp%	Act%
96.00	100	100
64.00	66.20	81.46#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

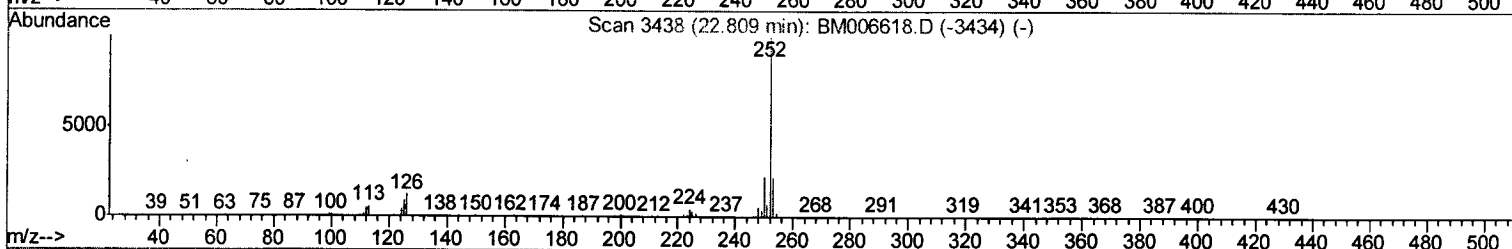
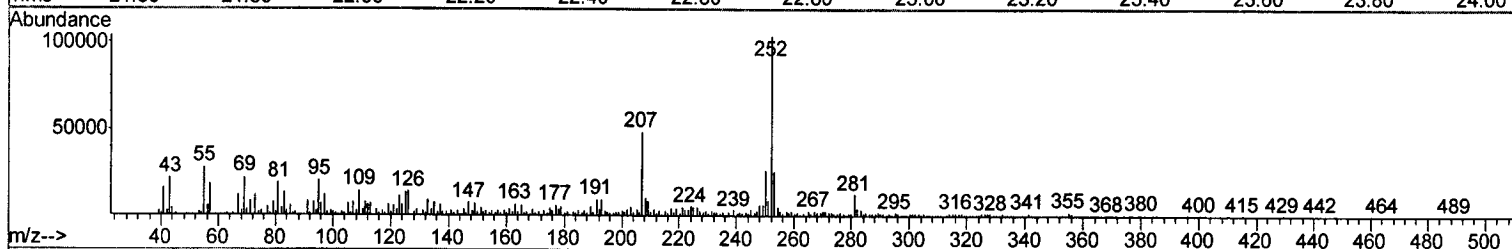
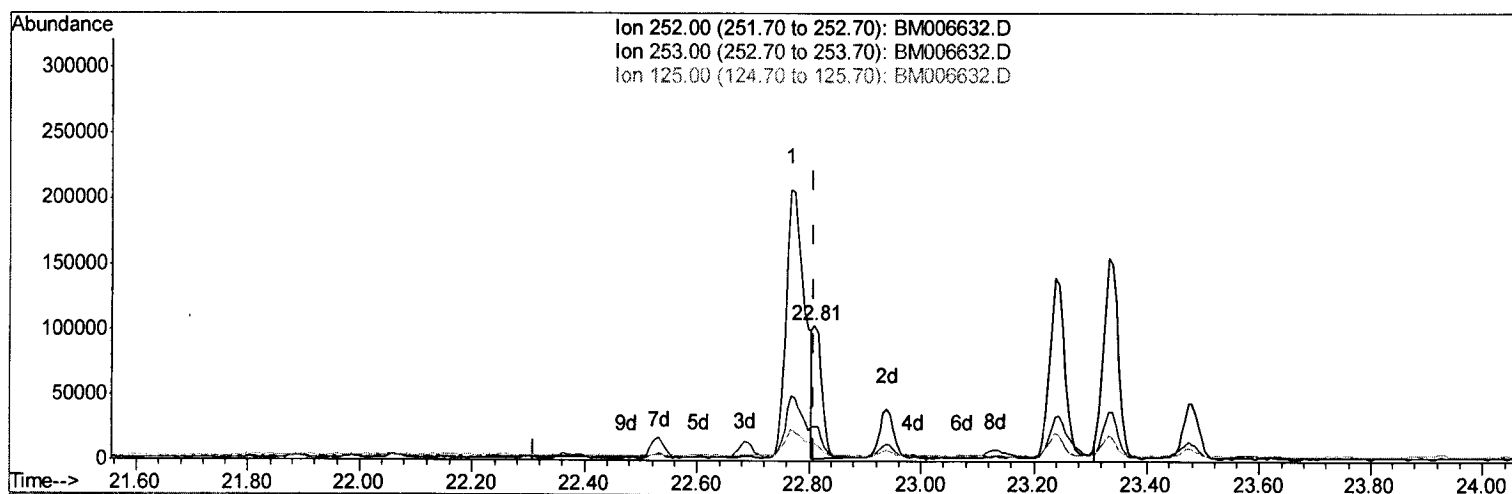
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TIC: BM006632.D

(36) Benzo(k)fluoranthene

22.809min (+0.000) 1.84ng/ul m

response 112525

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Ion	Exp%	Act%
252.00	100	100
253.00	22.30	25.23
125.00	9.60	13.52#
0.00	0.00	0.00

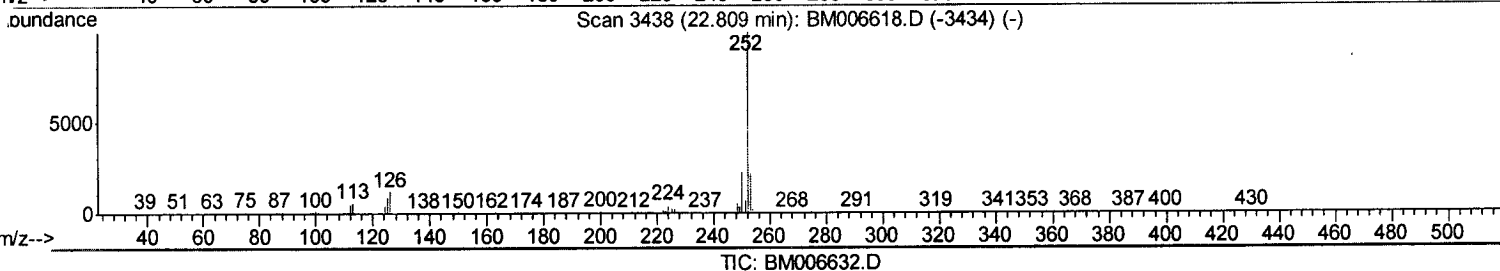
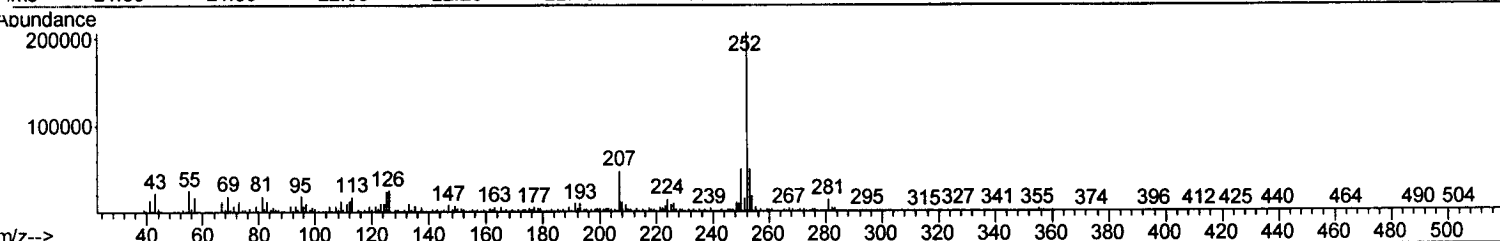
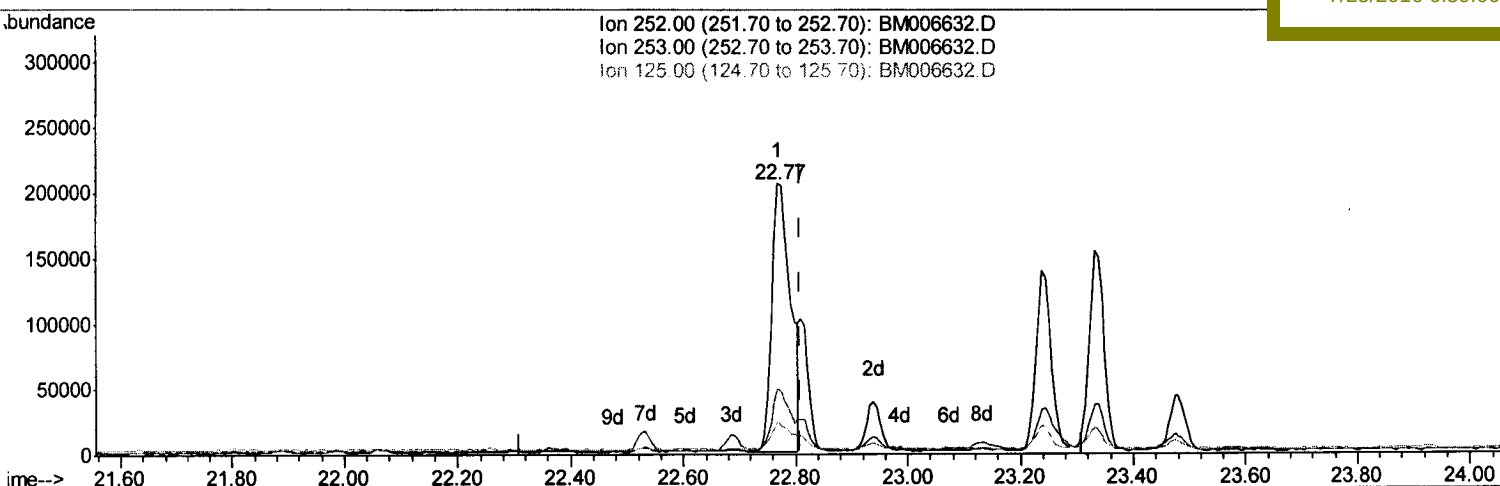
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Manual Integration

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(36) Benzo(k)fluoranthene

22.768min (-0.041) 7.77ng/ul

response 475450

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	23.96
125.00	9.60	11.74#
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	7.62	152	182825	20.00	ng/ul	0.00
7) Naphthalene-d8	10.40	136	803993	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.27	164	450462	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.02	188	990354	20.00	ng/ul	0.00
29) Chrysene-d12	21.22	240	1023544	20.00	ng/ul	0.00
34) Perylene-d12	23.43	264	1055433	20.00	ng/ul	0.01

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.13	96	10211m	2.26	ng/uL	0.00
3) Phenol-d5	6.80	99	446945	28.73	ng/ul	0.00
4) Bis-(2-Chloroethyl) ether-d	6.96	67	273877	28.42	ng/ul	0.00
5) 2-Chlorophenol-d4	7.16	132	351409	28.27	ng/ul	0.00
6) 4-Methylphenol-d8	8.33	113	349948	29.01	ng/ul	0.00
8) Nitrobenzene-d5	8.77	128	171907	28.25	ng/ul	0.00
9) 2-Nitrophenol-d4	9.49	143	184516	28.51	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.03	165	335244	27.41	ng/ul	0.00
12) 4-Chloroaniline-d4	10.55	131	302739	22.48	ng/ul	0.01
16) Dimethylphthalate-d6	13.69	166	964473	27.31	ng/ul	0.00
17) Acenaphthylene-d8	13.96	160	1272205	28.22	ng/ul	0.00
20) 4-Nitrophenol-d4	14.50	143	163162	26.60	ng/ul	0.01
21) Fluorene-d10	15.27	176	873722	27.66	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.41	200	59747	11.62	ng/ul	0.00
26) Anthracene-d10	17.12	188	1294453	27.53	ng/ul	0.00
30) Pyrene-d10	19.43	212	1389614	29.80	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.29	264	1348600	27.78	ng/ul	0.01

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Target Compounds

						Qvalue
19) Acenaphthene	14.34	153	824842	26.58	ng/ul	95
25) Phenanthrene	17.06	178	229553	4.12	ng/ul	100
27) Anthracene	17.16	178	68248	1.19	ng/ul	98
28) Fluoranthene	19.09	202	467935	7.15	ng/ul	99
31) Pyrene	19.46	202	2107194	34.00	ng/ul	97
32) Benzo(a)anthracene	21.21	228	236443	3.81	ng/ul	94
33) Chrysene	21.26	228	283095	4.95	ng/ul	97
35) Benzo(b)fluoranthene	22.77	252	475450	7.51	ng/ul#	96
36) Benzo(k)fluoranthene	22.81	252	112525m	1.84	ng/ul	94
38) Benzo(a)pyrene	23.33	252	280099	4.55	ng/ul#	94
39) Indeno(1,2,3-cd)pyrene	25.64	276	250697	3.52	ng/ul	99
40) Dibenzo(a,h)anthracene	25.64	278	70048	1.19	ng/ul#	86
41) Benzo(g,h,i)perylene	26.32	276	251949	4.12	ng/ul#	96

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