

Quantitation Report (QT Reviewed)

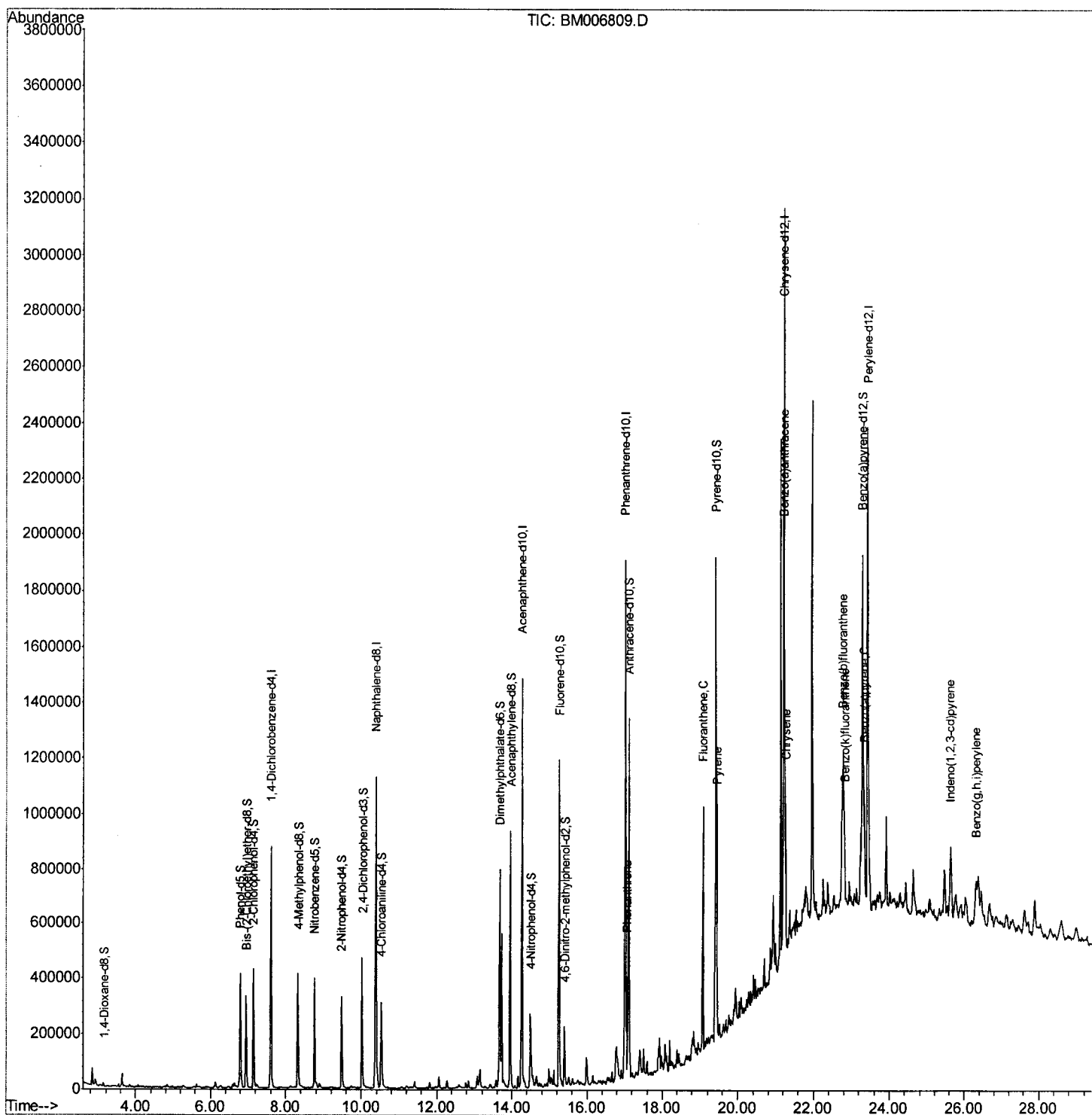
Data Path : Z:\HPCHEM1\BNA_M\Data\BM073116\
 Data File : BM006809.D
 Acq On : 01 Aug 2016 03:46
 Operator : UM/SJ
 Sample : H4190-09
 Misc :
 ALS Vial : 77 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 D9ZN8

Manual Integrations
 APPROVED

sohil
 8/1/2016 7:08:23 PM

Quant Time: Aug 01 08:09:47 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)

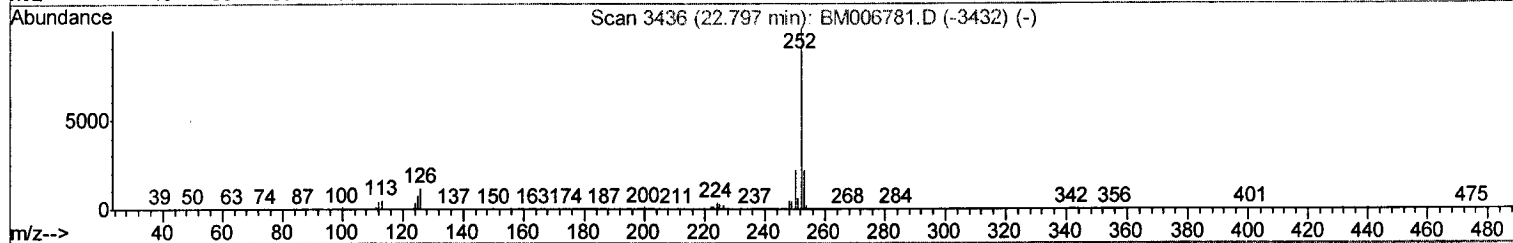
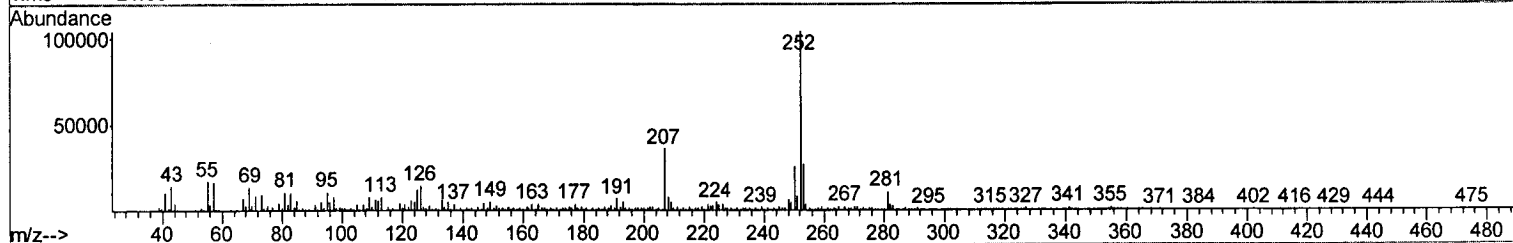
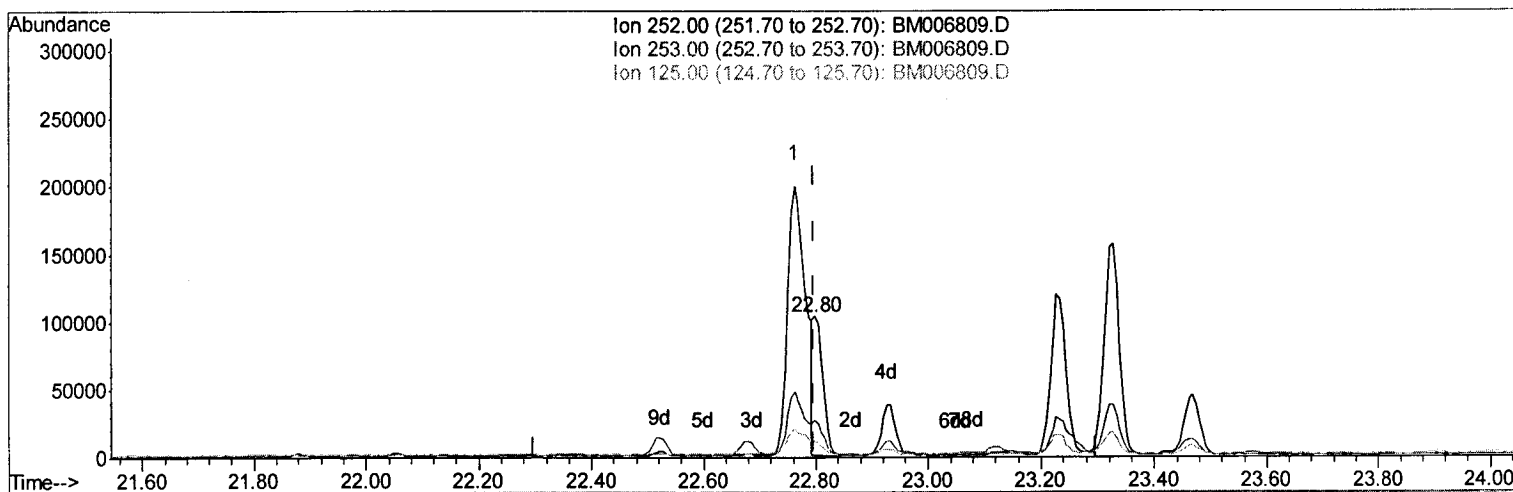
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TIC: BM006809.D

(36) Benzo(k)fluoranthene

22.797min (+0.000) 1.66ng/ul m

response 123012

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	26.04
125.00	9.60	11.59#
0.00	0.00	0.00

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Quantitation Report (Qedit)

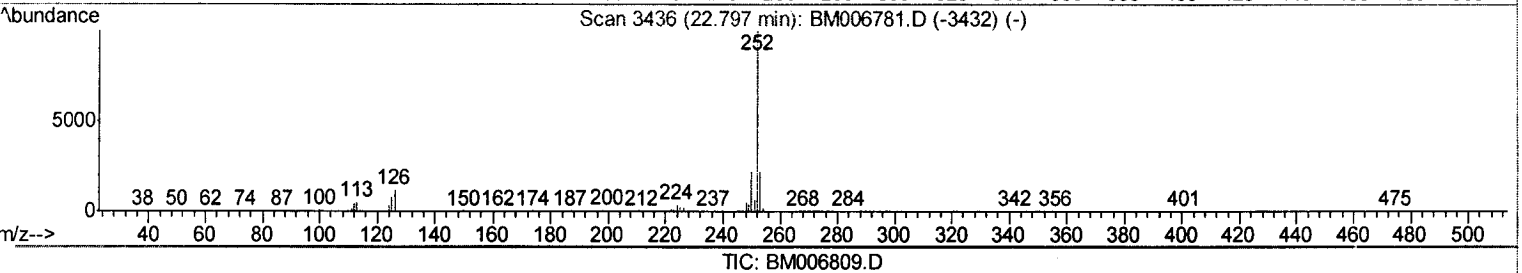
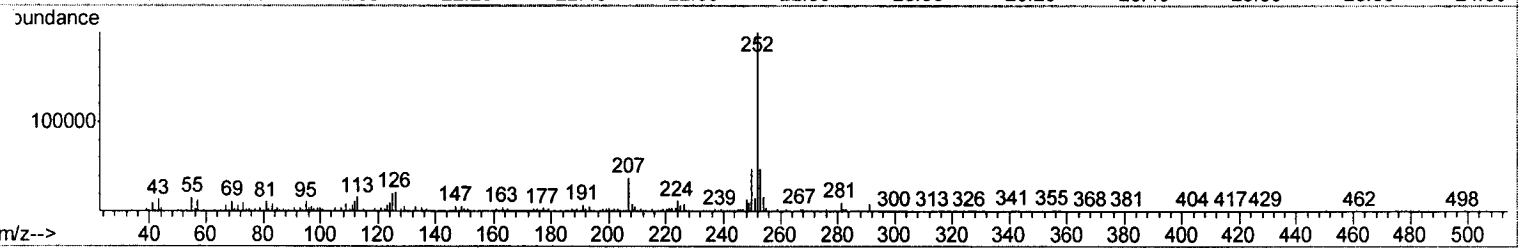
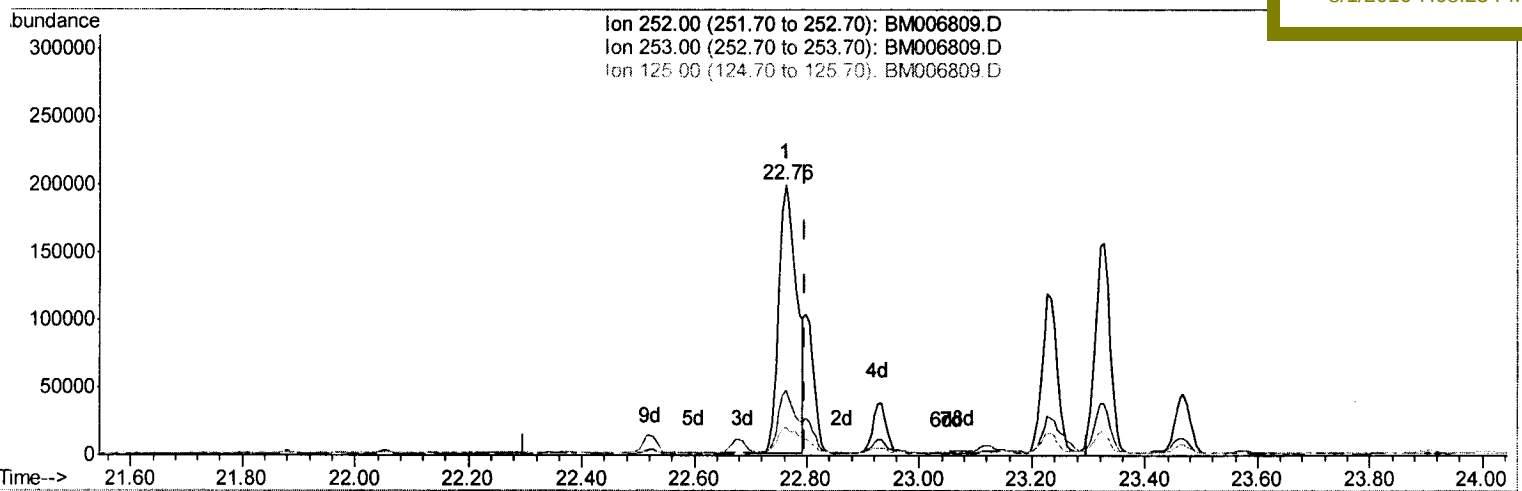
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(36) Benzo(k)fluoranthene

22.762min (-0.035) 5.99ng/ul

response 443123

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	24.30
125.00	9.60	10.68
0.00	0.00	0.00

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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	219704	20.00	ng/ul	0.00
7) Naphthalene-d8	10.38	136	931297	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.26	164	503492	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.02	188	1116604	20.00	ng/ul	0.00
29) Chrysene-d12	21.22	240	1204423	20.00	ng/ul	0.00
34) Perylene-d12	23.42	264	1275572	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.13	96	6045	1.11	ng/uL	0.00
3) Phenol-d5	6.79	99	228379	12.22	ng/ul	0.00
4) Bis-(2-Chloroethyl) ether-d	6.95	67	151767	13.11	ng/ul	0.00
5) 2-Chlorophenol-d4	7.14	132	190420	12.75	ng/ul	0.00
6) 4-Methylphenol-d8	8.32	113	158304	10.92	ng/ul	0.00
8) Nitrobenzene-d5	8.76	128	94359	13.39	ng/ul	0.00
9) 2-Nitrophenol-d4	9.48	143	100281	13.38	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.02	165	173397	12.24	ng/ul	0.00
12) 4-Chloroaniline-d4	10.53	131	175867	11.27	ng/ul	0.00
16) Dimethylphthalate-d6	13.67	166	523143	13.25	ng/ul	0.00
17) Acenaphthylene-d8	13.95	160	681576	13.53	ng/ul	0.00
20) 4-Nitrophenol-d4	14.50	143	65796	9.60	ng/ul	0.02
21) Fluorene-d10	15.26	176	472467	13.38	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.41	200	41052	7.08	ng/ul	0.01
26) Anthracene-d10	17.11	188	728501	13.74	ng/ul	0.00
30) Pyrene-d10	19.42	212	821218	14.96	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.28	264	818253	13.95	ng/ul	0.00

Target Compounds

						Qvalue
25) Phenanthrene	17.06	178	220928	3.52	ng/ul	99
28) Fluoranthene	19.08	202	520369	7.05	ng/ul	99
31) Pyrene	19.44	202	441164	6.05	ng/ul	98
32) Benzo(a)anthracene	21.20	228	287078	3.93	ng/ul	99
33) Chrysene	21.25	228	280602	4.17	ng/ul	97
35) Benzo(b)fluoranthene	22.76	252	443123	5.80	ng/ul	97
36) Benzo(k)fluoranthene	22.80	252	123012m	1.66	ng/ul	97
38) Benzo(a)pyrene	23.33	252	288451	3.88	ng/ul	95
39) Indeno(1,2,3-cd)pyrene	25.62	276	228179	2.65	ng/ul	99
41) Benzo(g,h,i)perylene	26.30	276	205522	2.78	ng/ul#	95

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