

Quantitation Report (QT Reviewed)

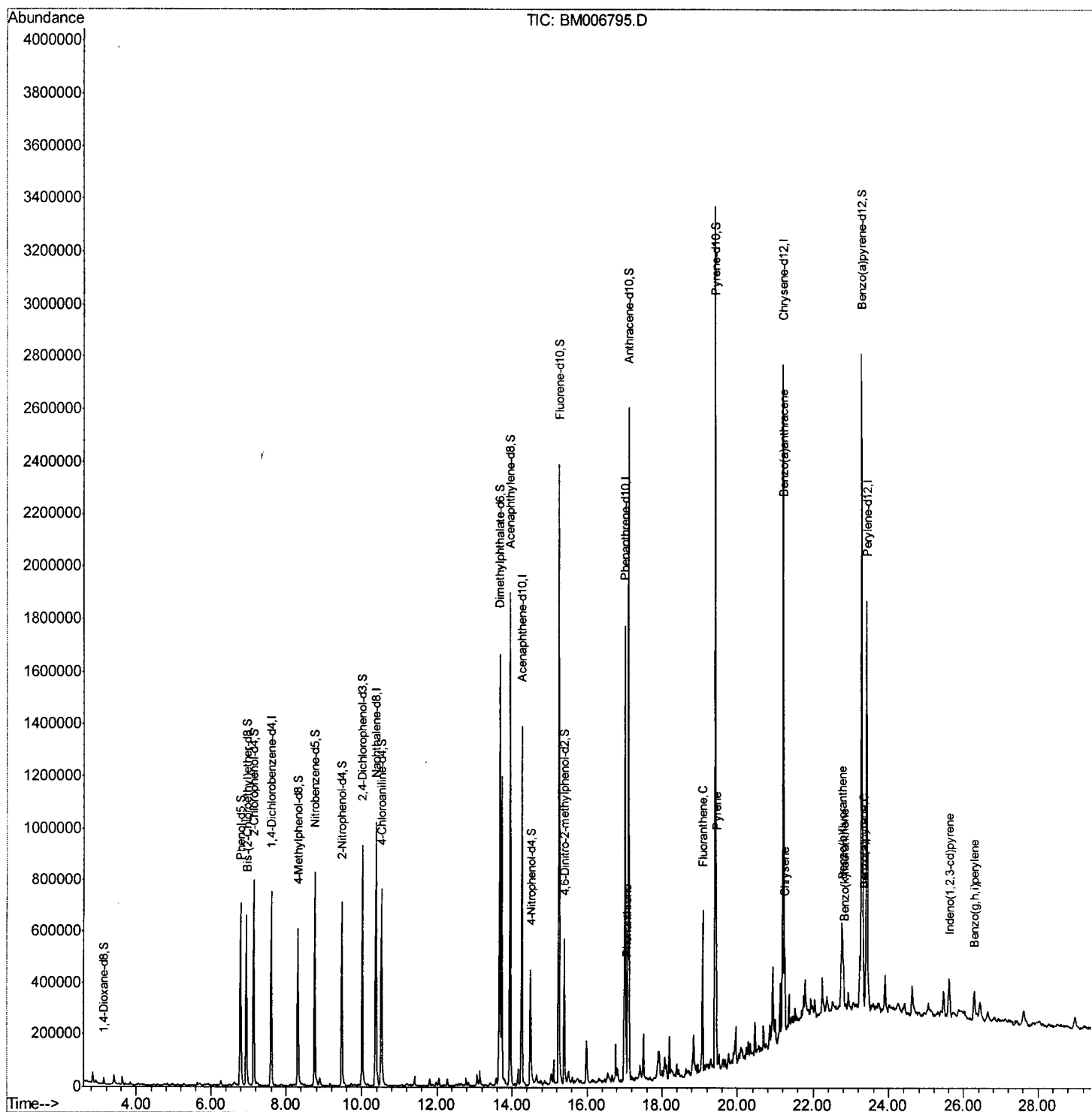
Data Path : Z:\HPCHEM1\BNA_M\Data\BM073116\
 Data File : BM006795.D
 Acq On : 31 Jul 2016 18:36
 Operator : UM/SJ
 Sample : H4189-03
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9ZM2

Manual Integrations
 APPROVED

sohil
 8/1/2016 7:07:19 PM

Quant Time: Aug 01 07:35:35 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)

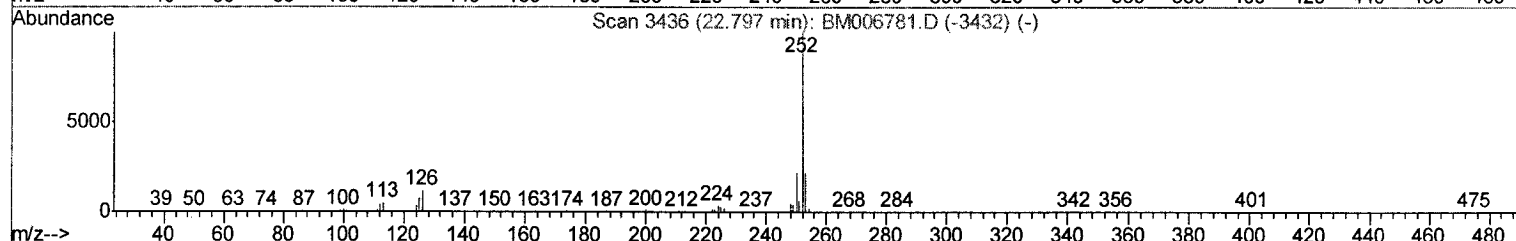
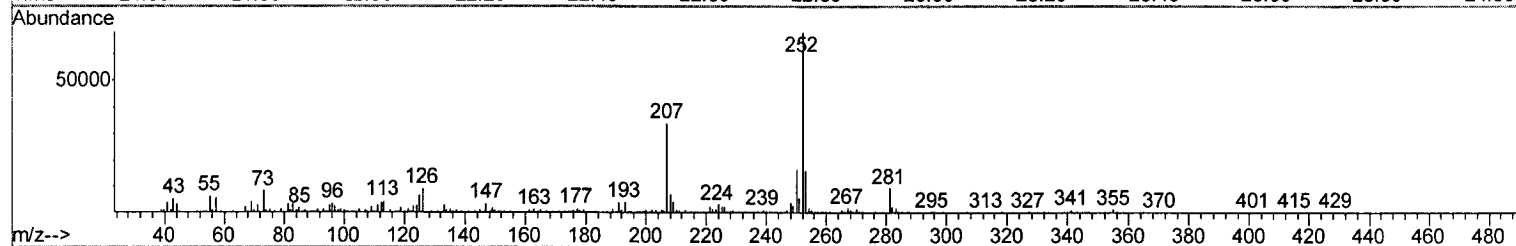
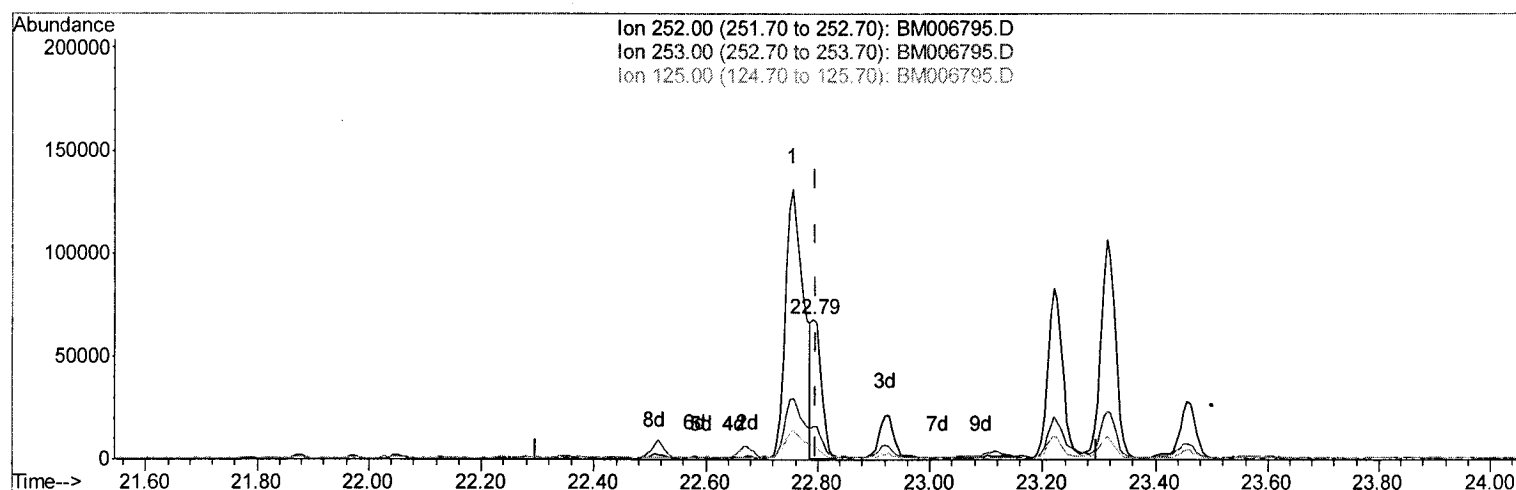
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TIC: BM006795.D

(36) Benzo(k)fluoranthene

22.791min (-0.006) 1.18ng/ul m U.M
 response 81554 08/02/16

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	23.90
125.00	9.60	9.96
0.00	0.00	0.00

Quantitation Report (Qedit)

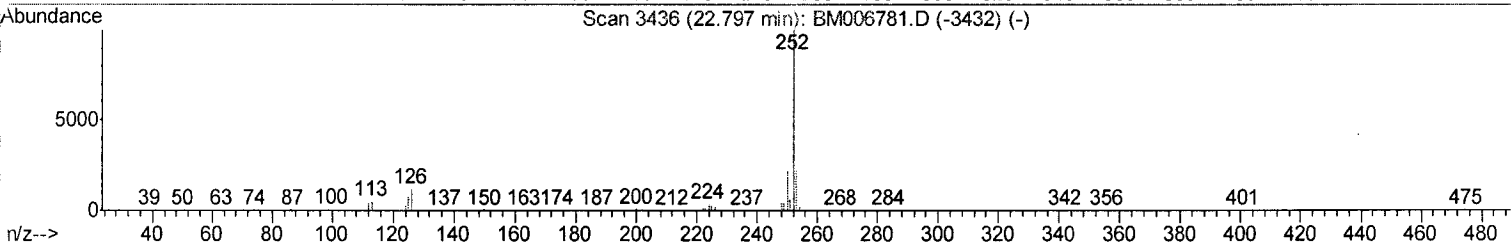
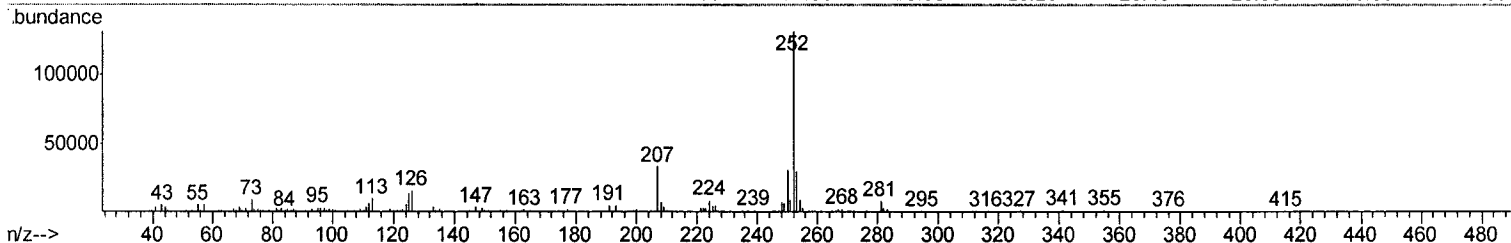
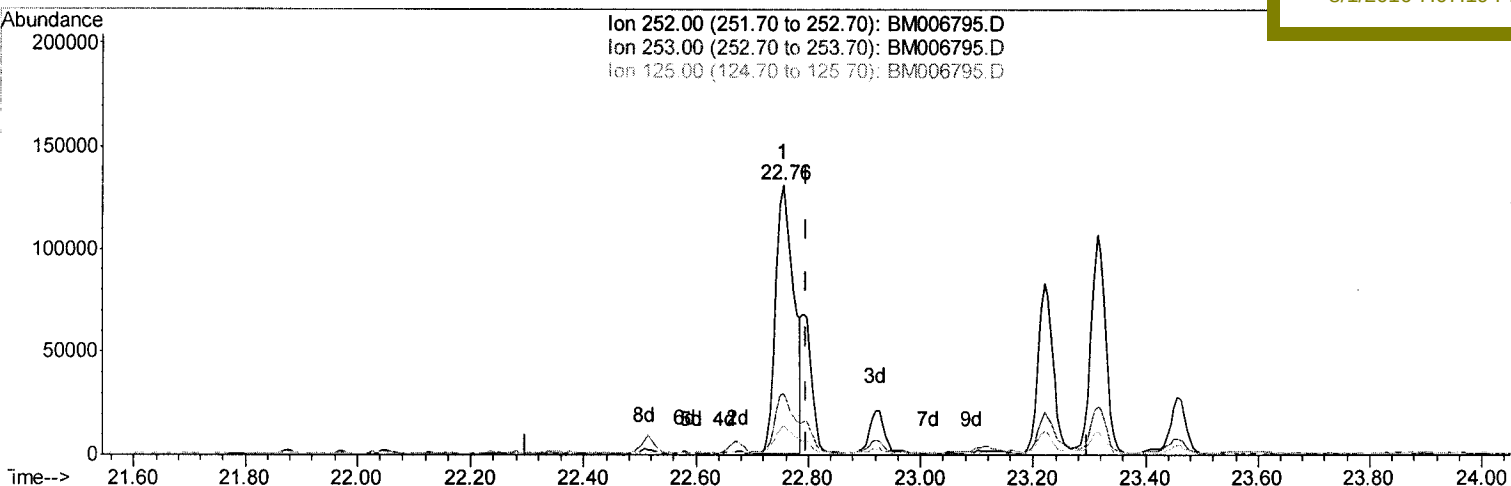
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Instrument :
 BNA_M
 ClientSampled :
 D9ZM2

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Manual Integrations
 APPROVED

sohil
 8/1/2016 7:07:19 PM



TIC: BM006795.D

(36) Benzo(k)fluoranthene

22.756min (-0.041) 4.31ng/ul

response 296620

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	22.81
125.00	9.60	10.84
0.00	0.00	0.00

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 Operator : UM/SJ
 Sample : H4189-03
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9ZM2

Manual Integrations
 APPROVED

sohil
 8/1/2016 7:07:19 PM

Quant Time: Aug 01 07:35:35 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	188565	20.00	ng/ul	0.00
7) Naphthalene-d8	10.38	136	835208	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.26	164	462510	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.01	188	1034630	20.00	ng/ul	0.00
29) Chrysene-d12	21.21	240	1157819	20.00	ng/ul	0.00
34) Perylene-d12	23.41	264	1188429	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.13	96	10491	2.25	ng/uL	0.00
3) Phenol-d5	6.79	99	387783	24.17	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.95	67	295244	29.71	ng/ul	0.00
5) 2-Chlorophenol-d4	7.14	132	345243	26.93	ng/ul	0.00
6) 4-Methylphenol-d8	8.32	113	231041	18.57	ng/ul	0.00
8) Nitrobenzene-d5	8.76	128	183044	28.95	ng/ul	0.00
9) 2-Nitrophenol-d4	9.47	143	205689	30.59	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.01	165	343791	27.06	ng/ul	0.00
12) 4-Chloroaniline-d4	10.53	131	403266	28.82	ng/ul	0.00
16) Dimethylphthalate-d6	13.67	166	1084436	29.90	ng/ul	0.00
17) Acenaphthylene-d8	13.94	160	1398055	30.20	ng/ul	0.00
20) 4-Nitrophenol-d4	14.49	143	118934	18.88	ng/ul	0.00
21) Fluorene-d10	15.26	176	965187	29.76	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.40	200	111089	20.69	ng/ul	0.00
26) Anthracene-d10	17.11	188	1453721	29.59	ng/ul	0.00
30) Pyrene-d10	19.42	212	1665829	31.58	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.27	264	1711922	31.32	ng/ul	0.00

Target Compounds

						Qvalue
25) Phenanthrene	17.05	178	175238	3.01	ng/ul	99
28) Fluoranthene	19.07	202	355731	5.20	ng/ul	99
31) Pyrene	19.44	202	306491	4.37	ng/ul	98
32) Benzo(a)anthracene	21.20	228	193560	2.76	ng/ul	99
33) Chrysene	21.24	228	204797	3.17	ng/ul	97
35) Benzo(b)fluoranthene	22.76	252	296620	4.16	ng/ul	99
36) Benzo(k)fluoranthene	22.79	252	81554m	1.18	ng/ul	99
38) Benzo(a)pyrene	23.31	252	189663	2.74	ng/ul	99
39) Indeno(1,2,3-cd)pyrene	25.61	276	139615	1.74	ng/ul	96
41) Benzo(g,h,i)perylene	26.29	276	136861	1.99	ng/ul	97

U.M
 08/02/16

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