

Quantitation Report (QT Reviewed)

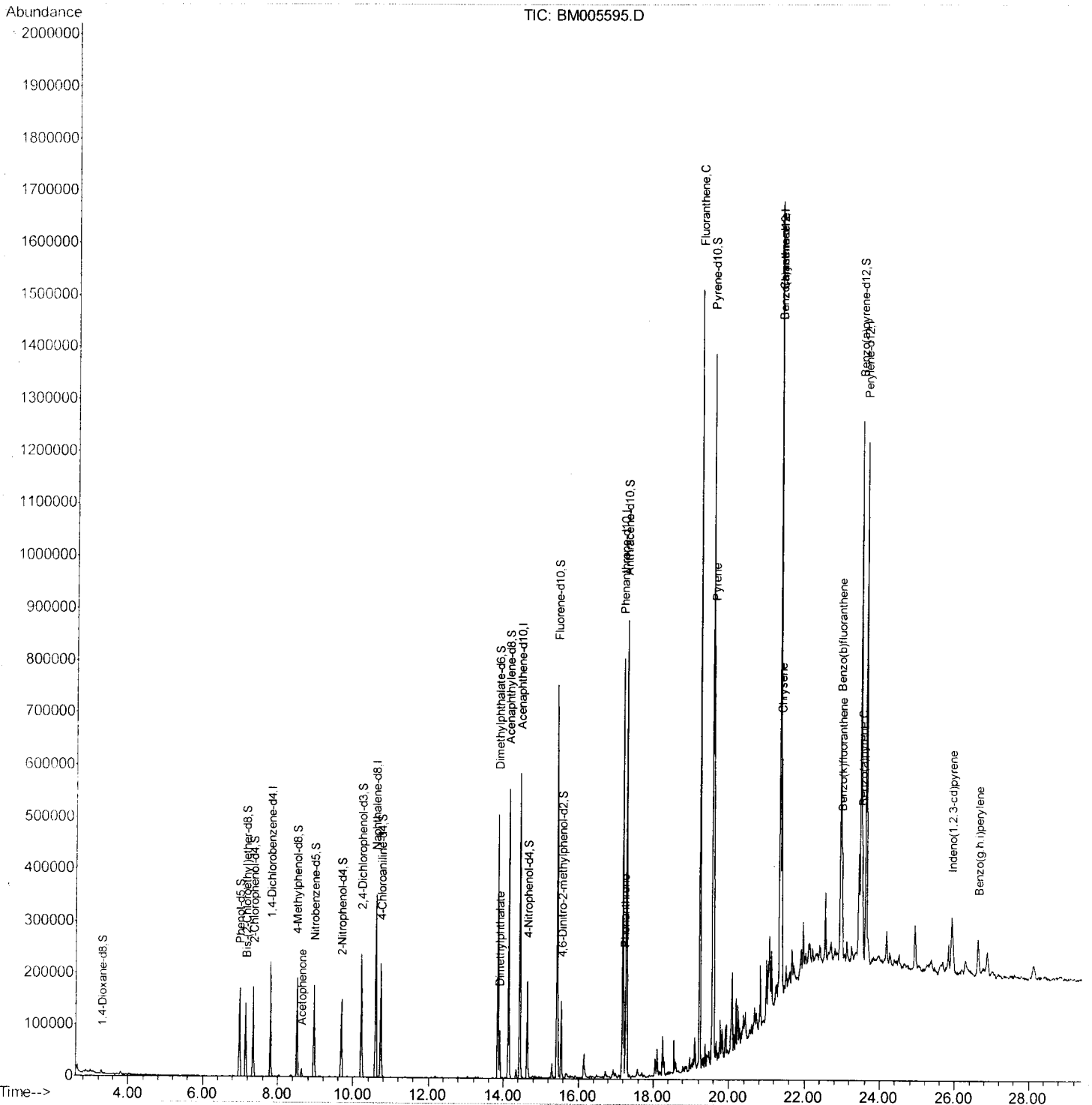
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052216\
 Data File : BM005595.D
 Acq On : 22 May 2016 20:15
 Operator : UM/SJ
 Sample : H3087-13
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHGK6

Manual Integrations
 APPROVED

umangi
 5/23/2016 8:35:47 PM

Quant Time: May 23 07:00:24 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM052016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon May 23 05:32:02 2016
 Response via : Initial Calibration



Quantitation Report (Qedit)

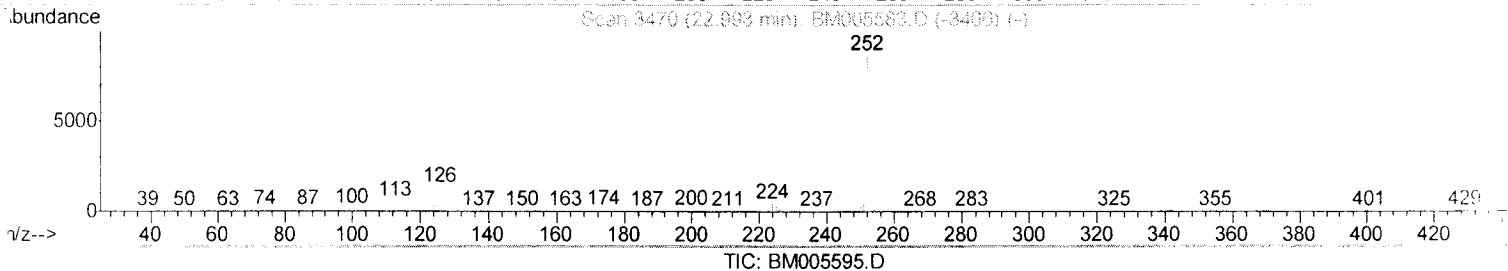
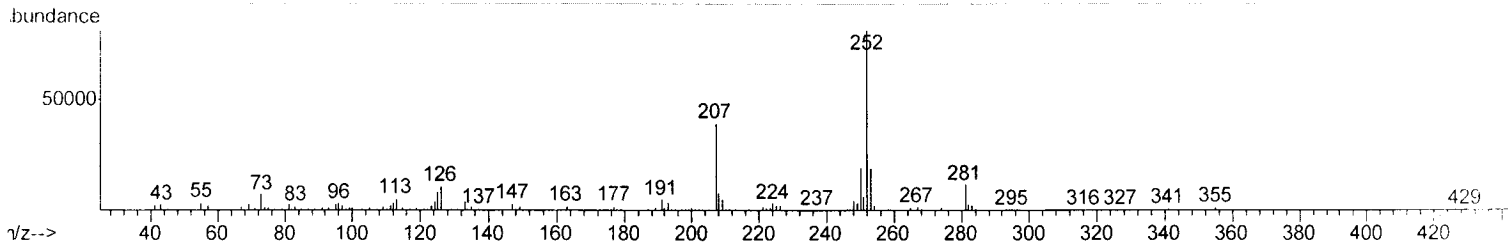
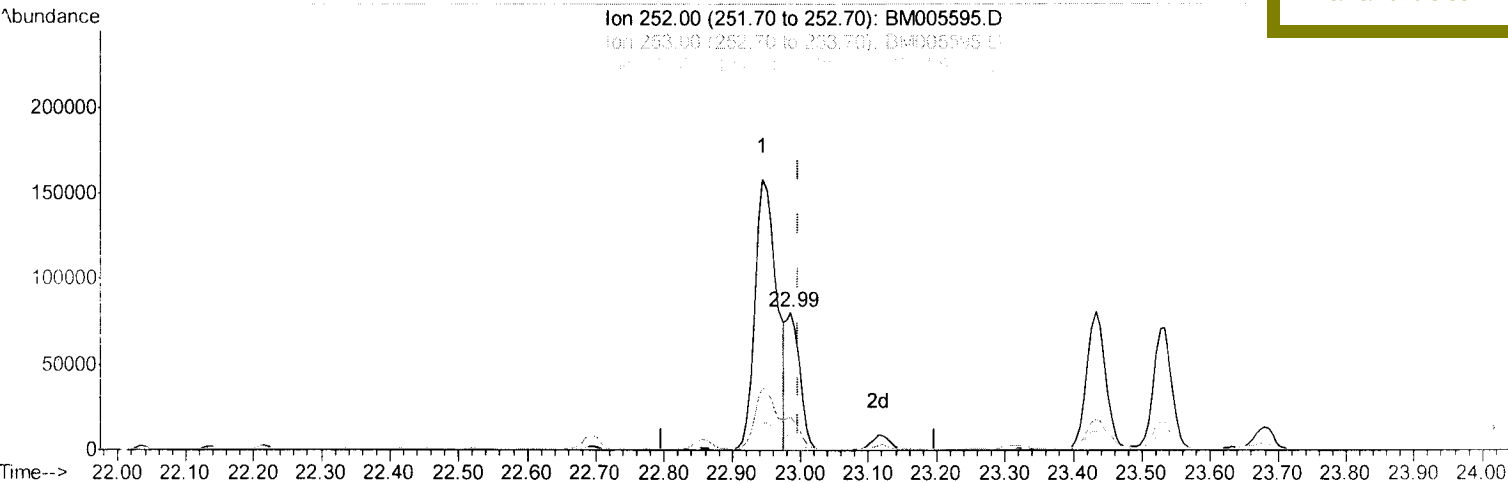
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(86) Benzo(k)fluoranthene

22.986min (-0.012) 2.80ng/ul m

response 117469

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	23.65
125.00	8.80	10.93#
0.00	0.00	0.00

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 05/25/16

Quantitation Report (Qedit)

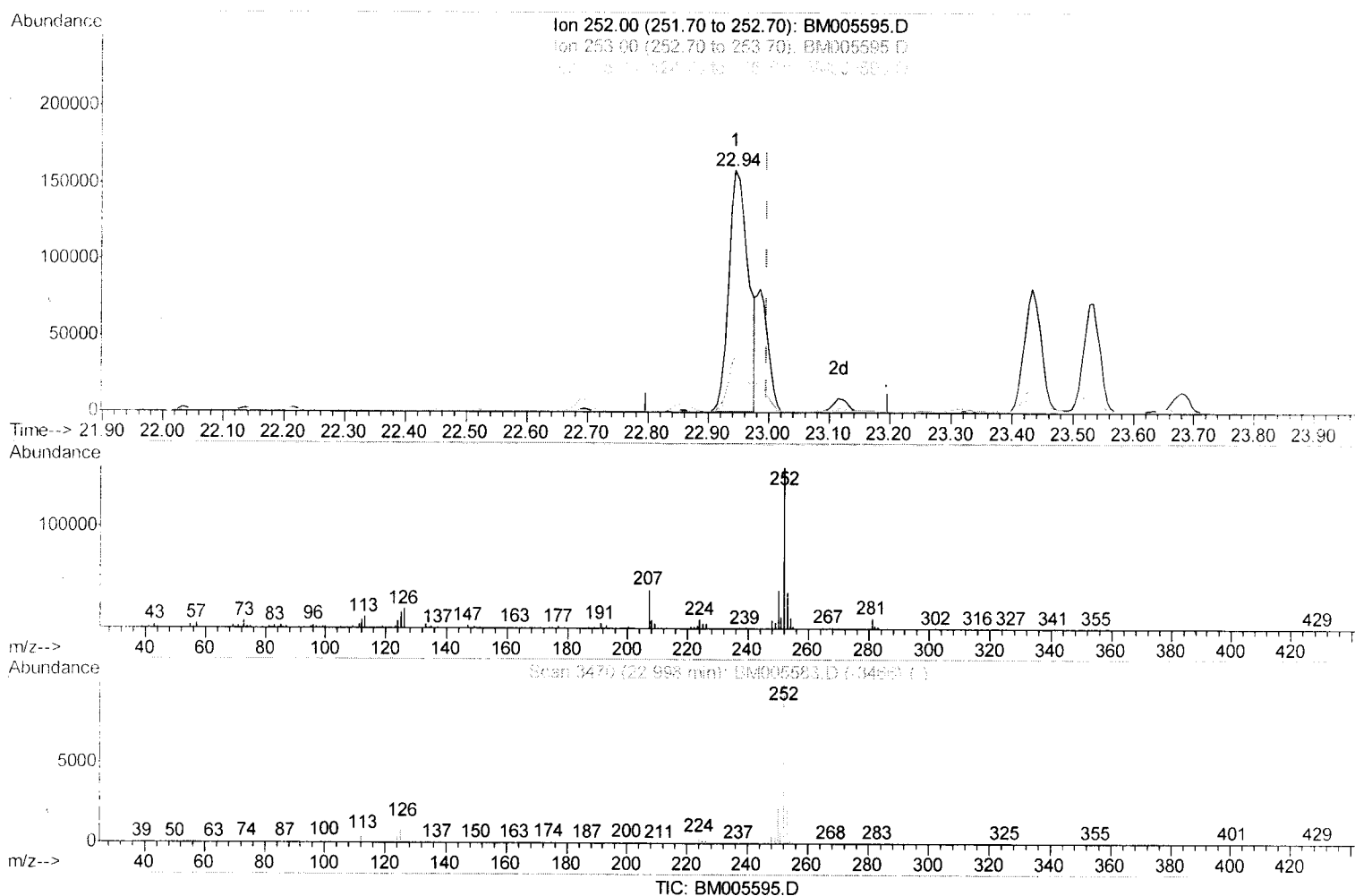
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(86) Benzo(k)fluoranthene

22.945min (-0.053) 8.25ng/ul

response 346310

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	23.11
125.00	8.80	10.44
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	7.81	152	59296	20.00	ng/ul	0.00
18) Naphthalene-d8	10.60	136	294852	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.45	164	190605	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.19	188	481575	20.00	ng/ul	0.00
75) Chrysene-d12	21.36	240	658472	20.00	ng/ul	0.00
83) Perylene-d12	23.63	264	752077	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	4276	3.16	ng/uL	0.00
5) Phenol-d5	6.98	99	99180	20.41	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.15	67	63888	22.68	ng/ul	0.00
9) 2-Chlorophenol-d4	7.35	132	82385	20.32	ng/ul	0.00
13) 4-Methylphenol-d8	8.51	113	73935	18.56	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	43506	19.30	ng/ul	0.00
22) 2-Nitrophenol-d4	9.69	143	49936	20.12	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.22	165	86889	18.59	ng/ul	0.00
29) 4-Chloroaniline-d4	10.73	131	119385	25.51	ng/ul	0.00
43) Dimethylphthalate-d6	13.85	166	324104	20.83	ng/ul	0.00
46) Acenaphthylene-d8	14.13	160	398053	20.46	ng/ul	0.00
51) 4-Nitrophenol-d4	14.64	143	44174	15.01	ng/ul	-0.01
57) Fluorene-d10	15.43	176	299632	22.13	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.55	200	31428	11.35	ng/ul	0.00
70) Anthracene-d10	17.28	188	488891	21.31	ng/ul	0.00
76) Pyrene-d10	19.57	212	604662	20.28	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.49	264	742835	21.15	ng/ul	-0.01

Target Compounds

					Qvalue
14) Acetophenone	8.63	105	8709	1.45 ng/ul	94
44) Dimethylphthalate	13.90	163	57431	3.74 ng/ul	99
69) Phenanthrene	17.23	178	101038	3.75 ng/ul	99
74) Fluoranthene	19.24	202	867092	28.62 ng/ul	97
77) Pyrene	19.60	202	494434	13.06 ng/ul	99
80) Benzo(a)anthracene	21.34	228	115710	2.99 ng/ul	97
82) Chrysene	21.39	228	269767	7.33 ng/ul	97
85) Benzo(b)fluoranthene	22.94	252	346310	7.81 ng/ul	96
86) Benzo(k)fluoranthene	22.99	252	117469m	2.80 ng/ul	
88) Benzo(a)pyrene	23.53	252	133700	3.11 ng/ul#	96
89) Indeno(1,2,3-cd)pyrene	25.92	276	104227	2.04 ng/ul	98
91) Benzo(g,h,i)perylene	26.63	276	83207	1.94 ng/ul	97

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 05/25/16

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