

Quantitation Report (QT Reviewed)

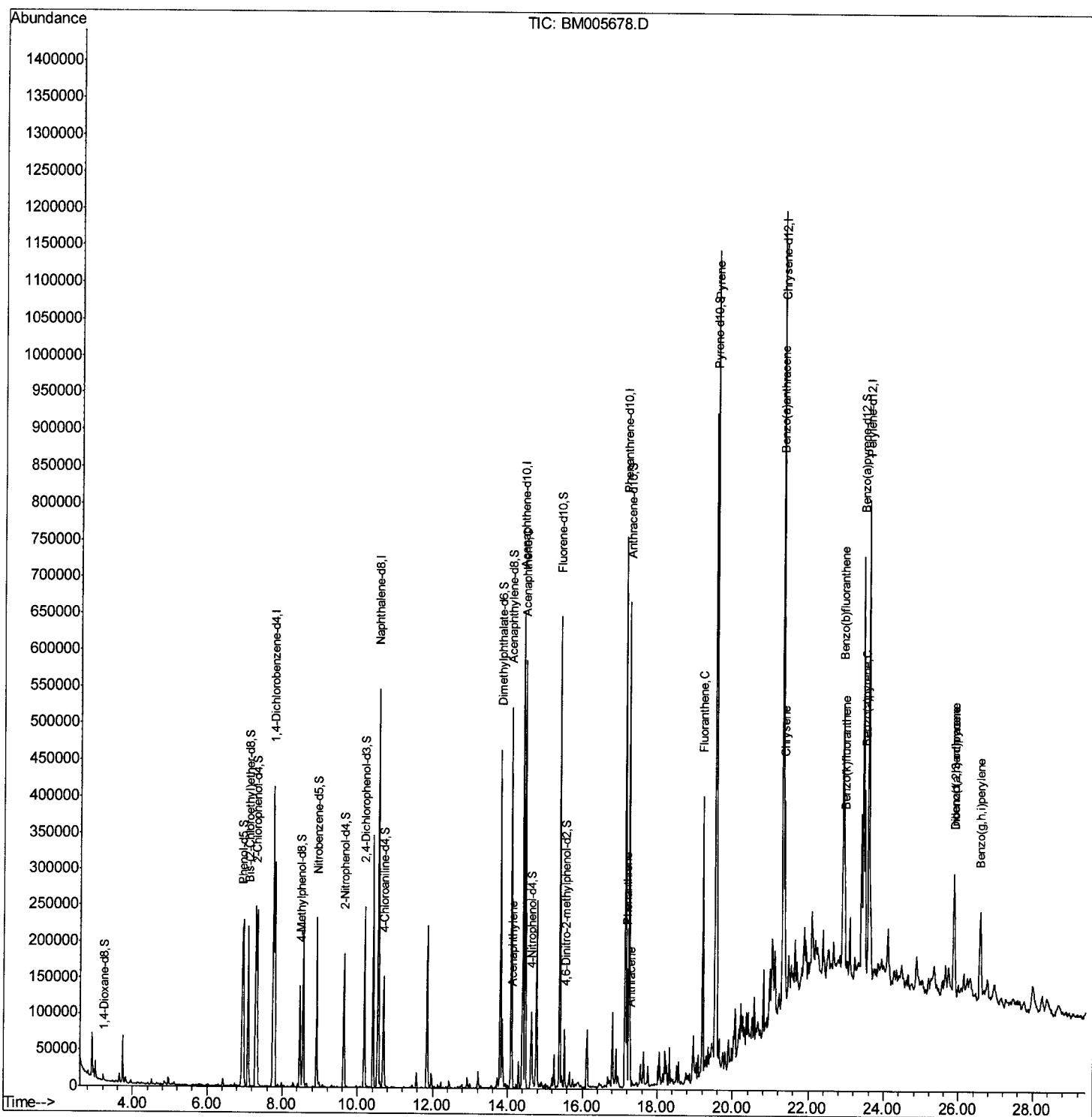
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060616\
 Data File : BM005678.D
 Acq On : 06 Jun 2016 20:14
 Operator : UM/SJ
 Sample : H3412-05MS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9XS5MS

Manual Integrations
 APPROVED

sohil
 6/7/2016 7:13:51 PM

Quant Time: Jun 07 04:59:10 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM060216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 07 04:25:20 2016
 Response via : Initial Calibration



Quantitation Report (Qedit)

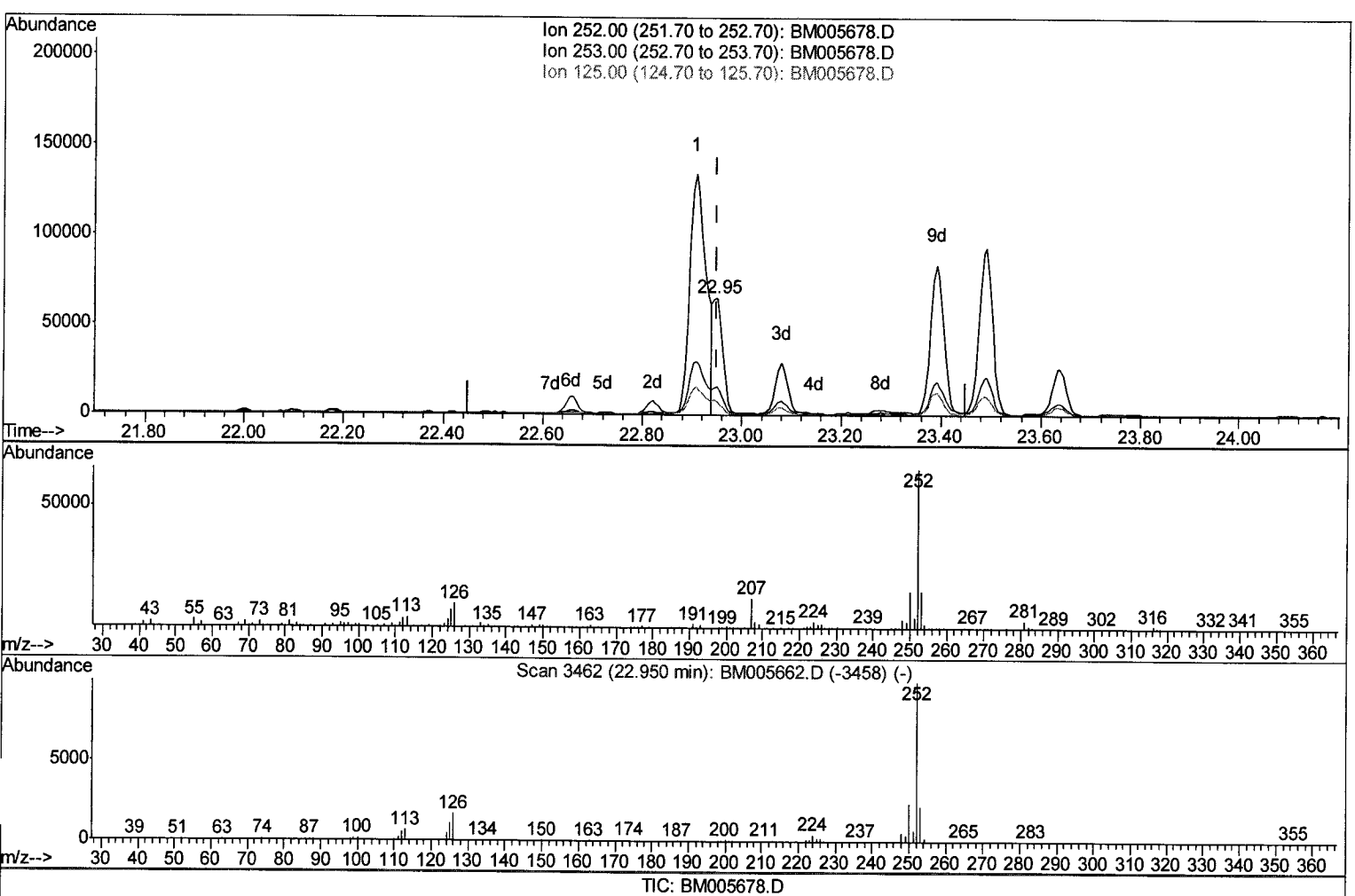
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(36) Benzo(k)fluoranthene

22.950min (+0.000) 3.37ng/ul m

response 87505

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	23.83
125.00	9.60	11.38
0.00	0.00	0.00

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 06/08/16

Quantitation Report (Qedit)

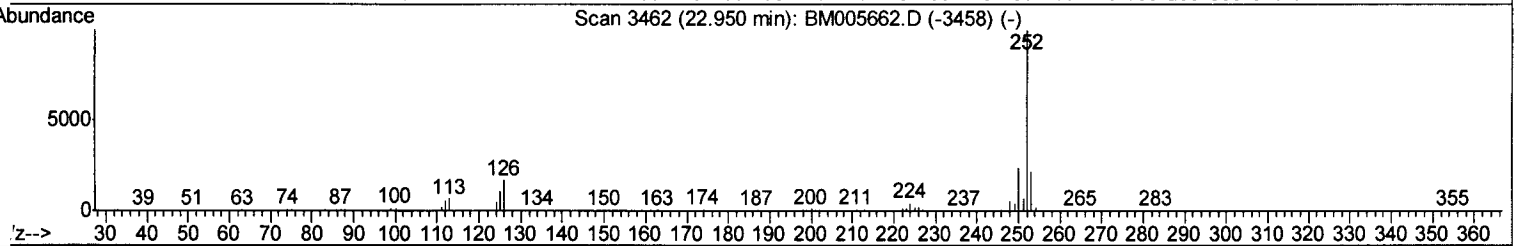
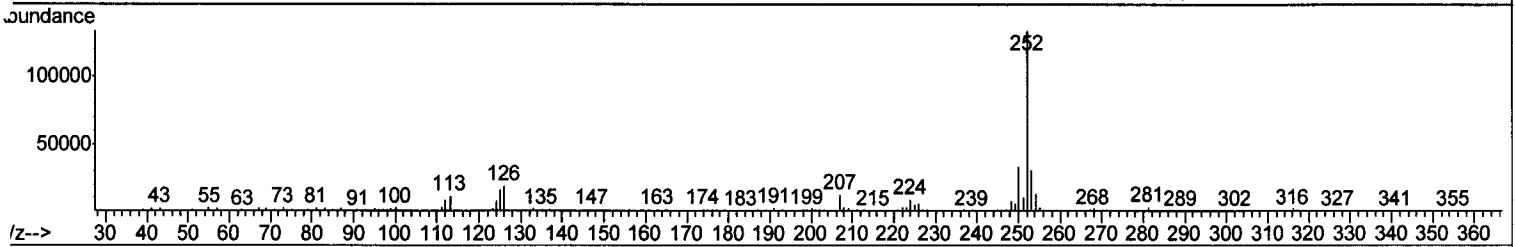
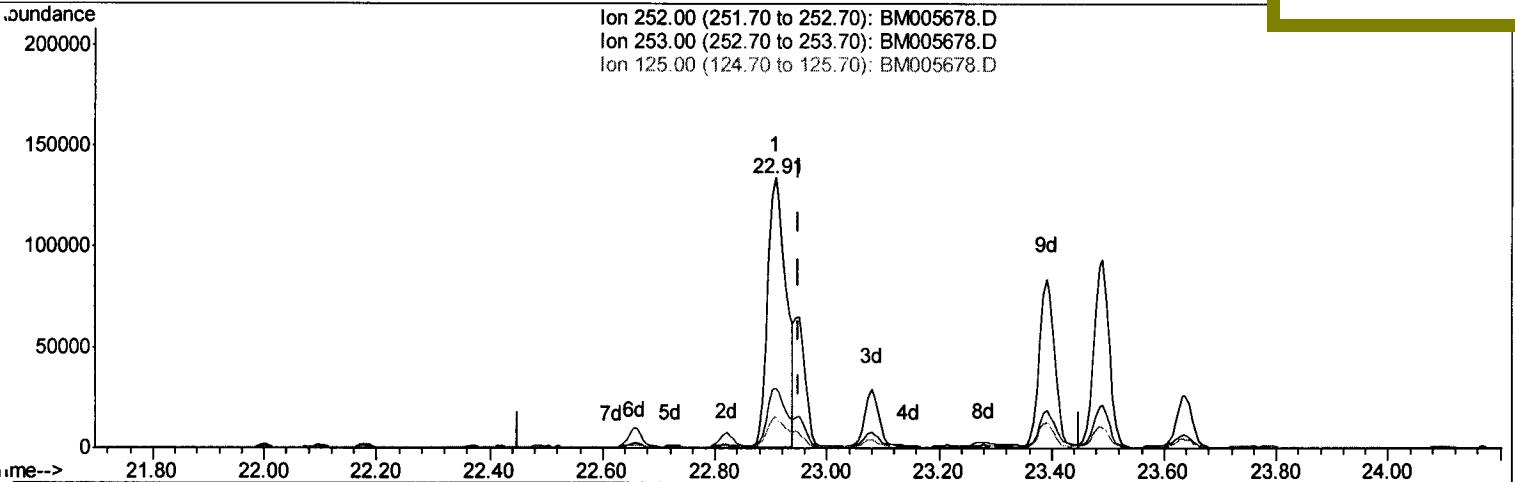
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TIC: BM005678.D

(36) Benzo(k)fluoranthene

22.909min (-0.041) 11.73ng/ul

response 304969

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	22.50
125.00	9.60	11.69#
0.00	0.00	0.00

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Manual Integrations
 APPROVED

sohil
 6/7/2016 7:13:51 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	109887	20.00	ng/ul	0.00
7) Naphthalene-d8	10.55	136	446732	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.40	164	213376	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.14	188	421378	20.00	ng/ul	0.00
29) Chrysene-d12	21.33	240	432434	20.00	ng/ul	0.00
34) Perylene-d12	23.59	264	459592	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.22	96	4161	1.59	ng/uL	0.00
3) Phenol-d5	6.94	99	123669	12.80	ng/ul	0.00
4) Bis-(2-Chloroethyl) ether-d	7.09	67	96730	16.45	ng/ul	0.00
5) 2-Chlorophenol-d4	7.29	132	116932	14.67	ng/ul	0.00
6) 4-Methylphenol-d8	8.47	113	54516	7.13	ng/ul	0.00
8) Nitrobenzene-d5	8.92	128	57475	16.75	ng/ul	0.00
9) 2-Nitrophenol-d4	9.64	143	59166	16.12	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.18	165	91642	14.05	ng/ul	0.00
12) 4-Chloroaniline-d4	10.69	131	87389	11.27	ng/ul	0.00
16) Dimethylphthalate-d6	13.81	166	290865	17.73	ng/ul	0.00
17) Acenaphthylene-d8	14.09	160	372805	17.47	ng/ul	0.00
20) 4-Nitrophenol-d4	14.63	143	17420	7.42	ng/ul	0.00
21) Fluorene-d10	15.39	176	251141	17.83	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	16239	8.18	ng/ul	0.00
26) Anthracene-d10	17.24	188	349070	17.44	ng/ul	0.00
30) Pyrene-d10	19.54	212	380575	17.51	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.44	264	374457	17.70	ng/ul	0.00

Target Compounds

					Qvalue	
18) Acenaphthylene	14.12	152	33420	1.51	ng/ul	100
19) Acenaphthene	14.46	153	233804	16.32	ng/ul	96
25) Phenanthrene	17.19	178	97932	4.21	ng/ul	99
27) Anthracene	17.28	178	31900	1.35	ng/ul	98
28) Fluoranthene	19.20	202	215500	9.06	ng/ul	99
31) Pyrene	19.57	202	632265	23.17	ng/ul	97
32) Benzo(a)anthracene	21.31	228	143620	5.66	ng/ul	96
33) Chrysene	21.36	228	152247	6.35	ng/ul	95
35) Benzo(b)fluoranthene	22.91	252	304969	11.00	ng/ul#	98
36) Benzo(k)fluoranthene	22.95	252	87505m	3.37	ng/ul	
38) Benzo(a)pyrene	23.49	252	172106	6.45	ng/ul	98
39) Indeno(1,2,3-cd)pyrene	25.87	276	147665	4.68	ng/ul	96
40) Dibenzo(a,h)anthracene	25.87	278	43388	1.64	ng/ul#	86
41) Benzo(g,h,i)perylene	26.57	276	144066	5.44	ng/ul#	94

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

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 6/8/16