

Quantitation Report (QT Reviewed)

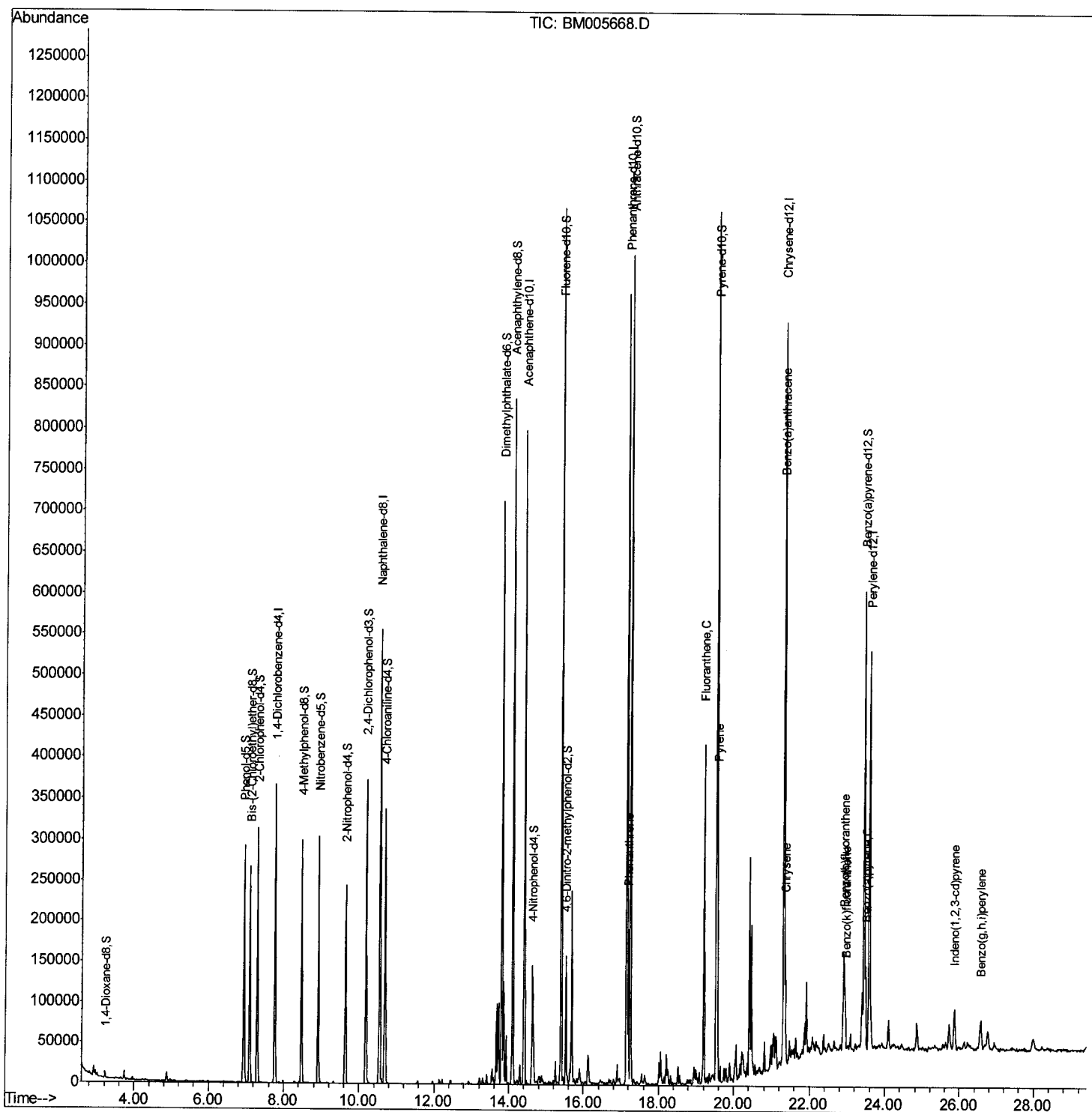
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060616\
 Data File : BM005668.D
 Acq On : 06 Jun 2016 14:02
 Operator : UM/SJ
 Sample : H3412-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9XT2

Manual Integrations
 APPROVED

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 6/7/2016 7:13:35 PM

Quant Time: Jun 07 04:38:46 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)

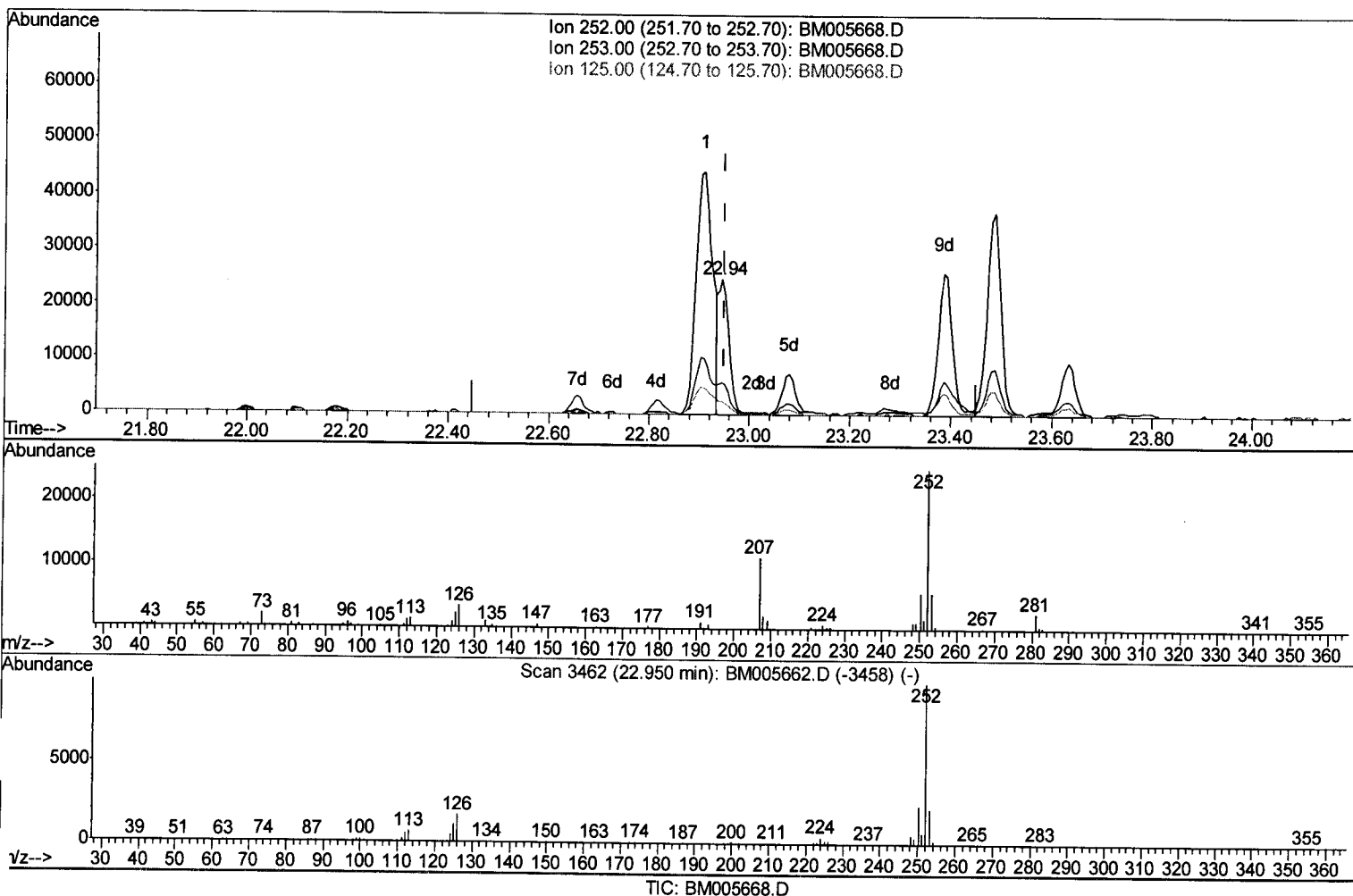
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(36) Benzo(k)fluoranthene

22.945min (-0.005) 1.85ng/ul m

response 36019

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	23.67
125.00	9.60	10.12
0.00	0.00	0.00

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Quantitation Report (Qedit)

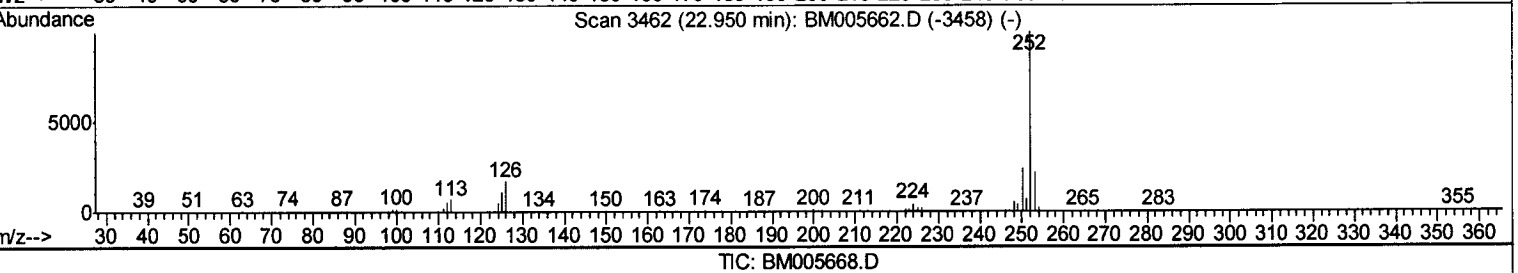
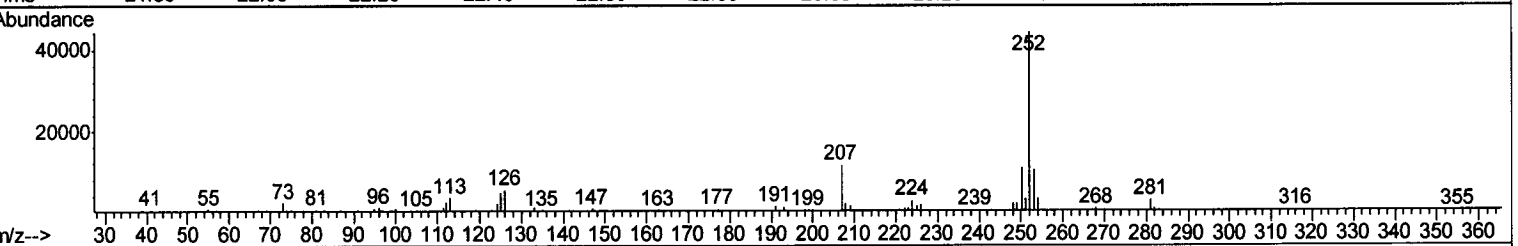
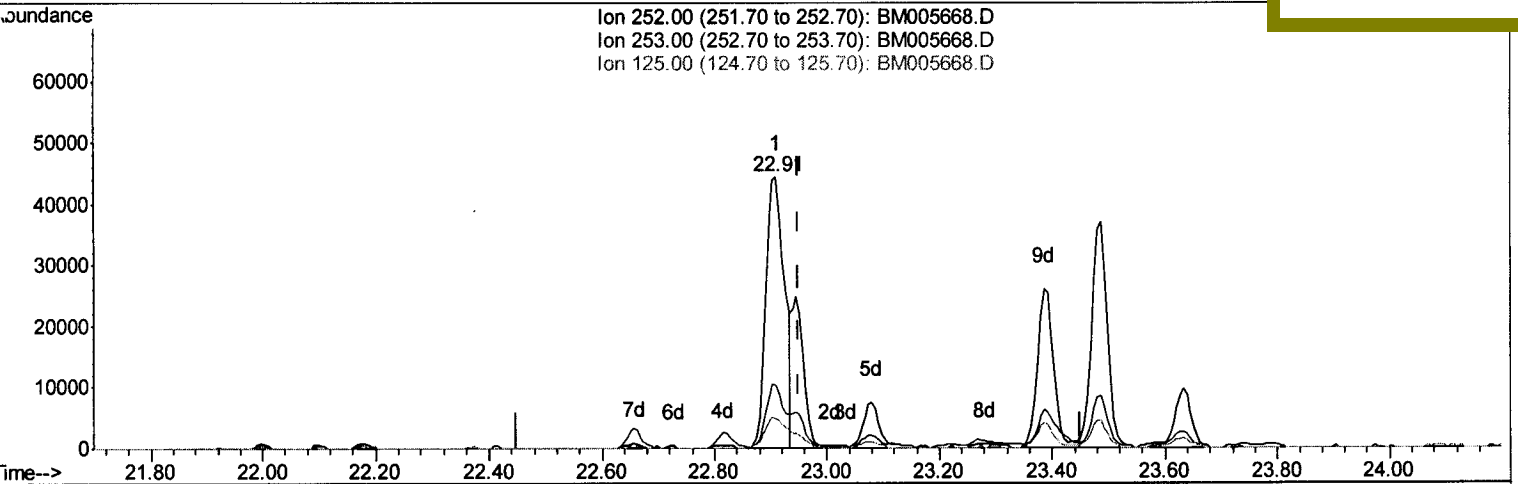
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(36) Benzo(k)fluoranthene

22.909min (-0.041) 5.22ng/ul

response 101480

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	23.03
125.00	9.60	10.96
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	7.76	152	97965	20.00	ng/ul	0.00
7) Naphthalene-d8	10.55	136	474243	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.40	164	269583	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.15	188	539746	20.00	ng/ul	0.00
29) Chrysene-d12	21.33	240	381038	20.00	ng/ul	0.00
34) Perylene-d12	23.59	264	343410	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.22	96	4340	1.86	ng/uL	0.00
3) Phenol-d5	6.93	99	174812	20.30	ng/ul	0.00
4) Bis-(2-Chloroethyl) ether-d	7.09	67	118792	22.67	ng/ul	0.00
5) 2-Chlorophenol-d4	7.29	132	147461	20.75	ng/ul	0.00
6) 4-Methylphenol-d8	8.47	113	117208	17.21	ng/ul	0.00
8) Nitrobenzene-d5	8.92	128	74561	20.47	ng/ul	0.00
9) 2-Nitrophenol-d4	9.64	143	81239	20.85	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.18	165	139771	20.18	ng/ul	0.00
12) 4-Chloroaniline-d4	10.69	131	189843	23.07	ng/ul	0.00
16) Dimethylphthalate-d6	13.81	166	467478	22.56	ng/ul	0.00
17) Acenaphthylene-d8	14.09	160	593533	22.02	ng/ul	0.00
20) 4-Nitrophenol-d4	14.62	143	42229	14.23	ng/ul	0.00
21) Fluorene-d10	15.39	176	408364	22.94	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	32483	12.78	ng/ul	0.00
26) Anthracene-d10	17.25	188	564661	22.03	ng/ul	0.00
30) Pyrene-d10	19.53	212	522055	27.26	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.44	264	358942	22.71	ng/ul	0.00

Target Compounds

						Qvalue
25) Phenanthrene	17.19	178	111171	3.73	ng/ul	98
28) Fluoranthene	19.20	202	242713	7.97	ng/ul	100
31) Pyrene	19.56	202	194017	8.07	ng/ul	97
32) Benzo(a)anthracene	21.31	228	91467	4.09	ng/ul	97
33) Chrysene	21.36	228	90016	4.26	ng/ul	96
35) Benzo(b)fluoranthene	22.91	252	101480	4.90	ng/ul	98
36) Benzo(k)fluoranthene	22.94	252	36019m	1.85	ng/ul	
38) Benzo(a)pyrene	23.49	252	71785	3.60	ng/ul	97
39) Indeno(1,2,3-cd)pyrene	25.86	276	48836	2.07	ng/ul	95
41) Benzo(g,h,i)perylene	26.57	276	44405	2.24	ng/ul#	94

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