

Quantitation Report (QT Reviewed)

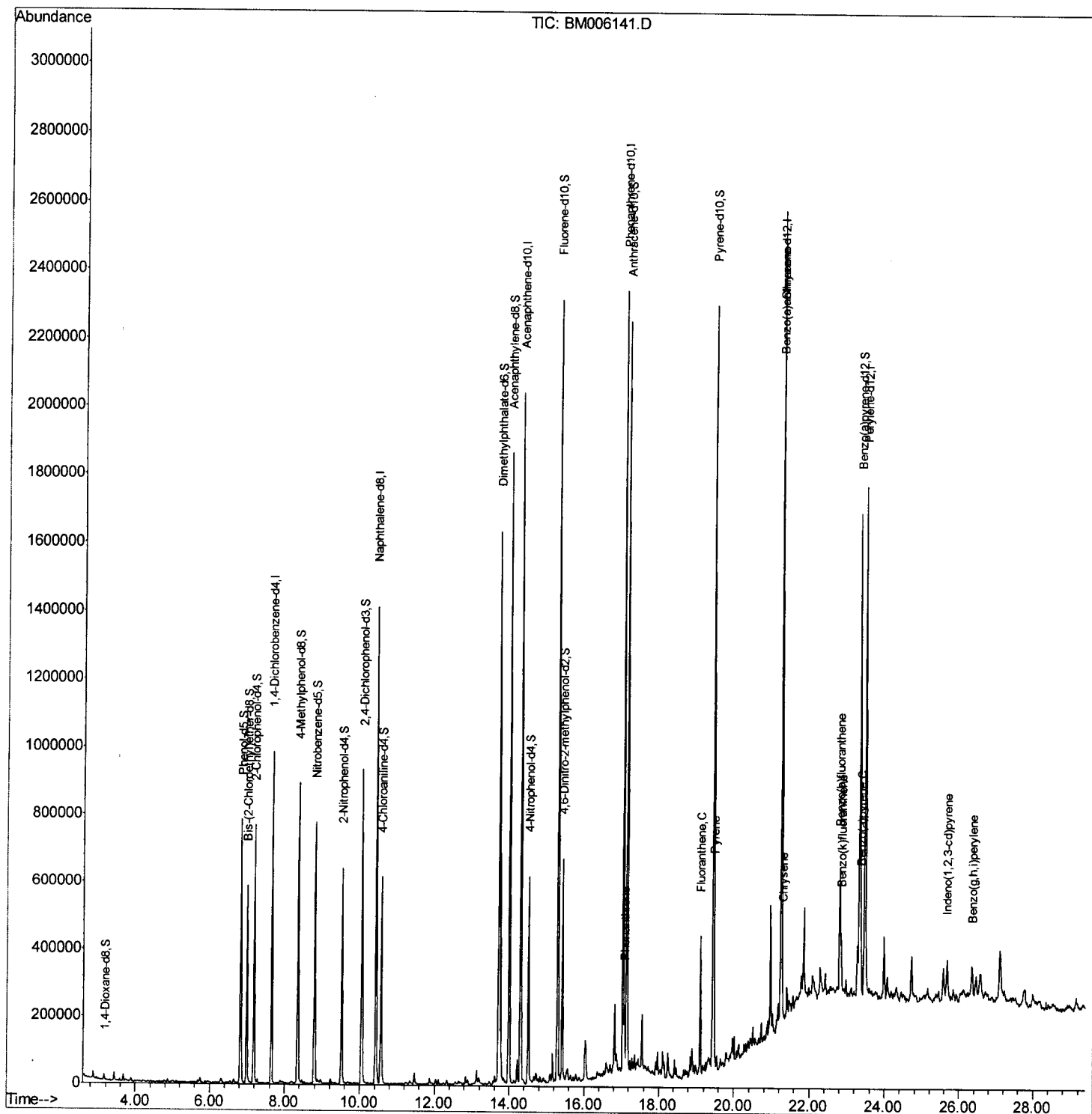
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062916\
 Data File : BM006141.D
 Acq On : 30 Jun 2016 00:39
 Operator : UM/SJ
 Sample : H3780-12
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9YG1

Manual Integration

Quant Time: Jun 30 04:07:57 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062816.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 29 00:57:09 2016
 Response via : Initial Calibration

sohil
 6/30/2016 7:44:22 PM



Quantitation Report (Qedit)

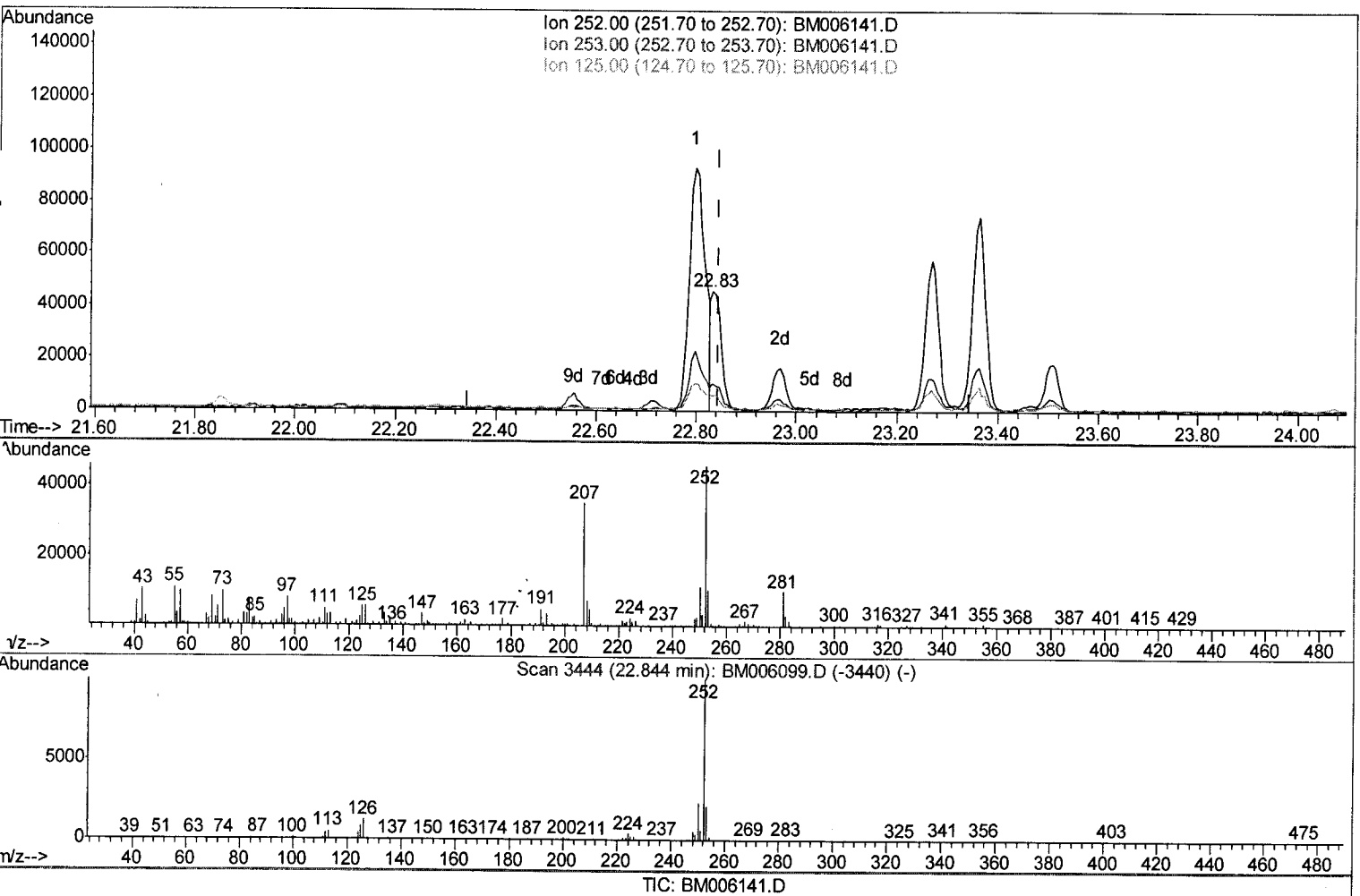
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(36) Benzo(k)fluoranthene

22.833min (-0.012) 1.17ng/ul m U.M.
 response 61441 07/02/16

Ion	Exp%	Act%
252.00	100	100
253.00	22.00	23.22
125.00	10.10	12.83#
0.00	0.00	0.00

Quantitation Report (Qedit)

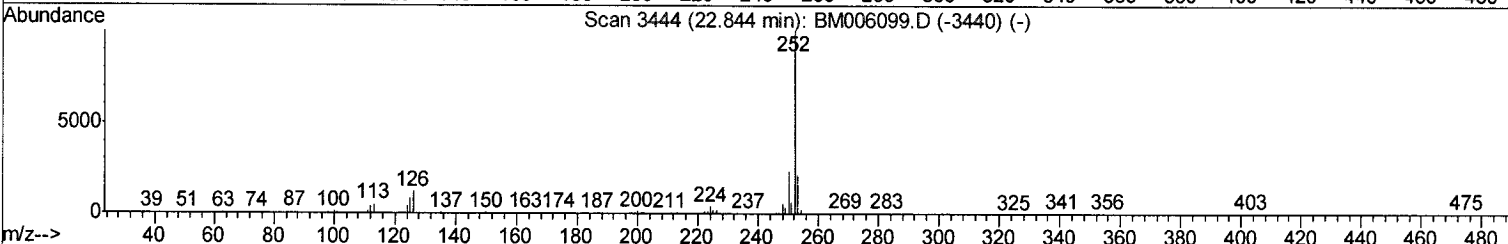
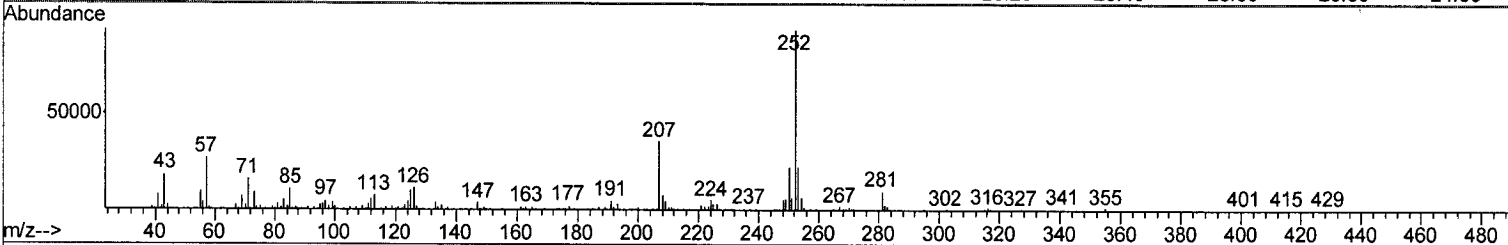
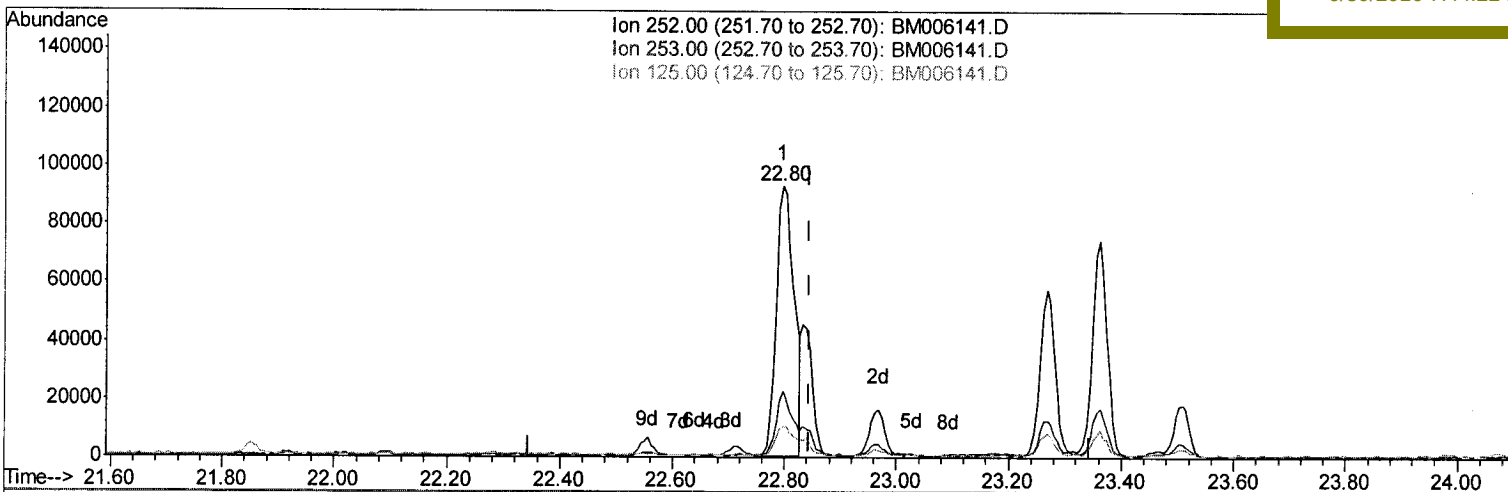
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 6/30/2016 7:44:22 PM



TIC: BM006141.D

(36) Benzo(k)fluoranthene

22.797min (-0.047) 3.90ng/ul

response 205355

Ion	Exp%	Act%
252.00	100	100
253.00	22.00	24.42
125.00	10.10	11.43
0.00	0.00	0.00

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Manual Integration

sohil
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	242835	20.00	ng/ul	0.00
7) Naphthalene-d8	10.43	136	1154013	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.30	164	667541	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.05	188	1331163	20.00	ng/ul	0.00
29) Chrysene-d12	21.24	240	1046752	20.00	ng/ul	0.00
34) Perylene-d12	23.46	264	1080515	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.16	96	8508	1.63	ng/uL	0.00
3) Phenol-d5	6.82	99	413816	20.86	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.99	67	263111	22.49	ng/ul	0.00
5) 2-Chlorophenol-d4	7.18	132	323135	21.22	ng/ul	0.00
6) 4-Methylphenol-d8	8.36	113	322436	20.19	ng/ul	0.00
8) Nitrobenzene-d5	8.80	128	171715	22.09	ng/ul	0.00
9) 2-Nitrophenol-d4	9.52	143	189027	21.25	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.05	165	341345	21.03	ng/ul	0.00
12) 4-Chloroaniline-d4	10.57	131	314330	24.50	ng/ul	0.00
16) Dimethylphthalate-d6	13.71	166	1078434	22.54	ng/ul	0.00
17) Acenaphthylene-d8	13.99	160	1342532	23.12	ng/ul	0.00
20) 4-Nitrophenol-d4	14.50	143	133509	14.71	ng/ul	0.00
21) Fluorene-d10	15.30	176	903140	22.12	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.42	200	122196	25.92	ng/ul	0.00
26) Anthracene-d10	17.14	188	1220296	22.85	ng/ul	0.00
30) Pyrene-d10	19.45	212	1139951	27.04	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.32	264	1006350	23.40	ng/ul	0.00

Target Compounds

					Qvalue	
25) Phenanthrene	17.09	178	114340	1.82	ng/ul	98
28) Fluoranthene	19.11	202	232349	3.01	ng/ul	98
31) Pyrene	19.47	202	196924	3.57	ng/ul	99
32) Benzo(a)anthracene	21.23	228	115686	2.13	ng/ul	97
33) Chrysene	21.28	228	128320	2.56	ng/ul	97
35) Benzo(b)fluoranthene	22.80	252	205355	3.66	ng/ul	94
36) Benzo(k)fluoranthene	22.83	252	61441m	1.17	ng/ul	97
38) Benzo(a)pyrene	23.36	252	133683	2.52	ng/ul#	93
39) Indeno(1,2,3-cd)pyrene	25.67	276	115172	2.09	ng/ul#	93
41) Benzo(g,h,i)perylene	26.36	276	115214	2.50	ng/ul	99

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