

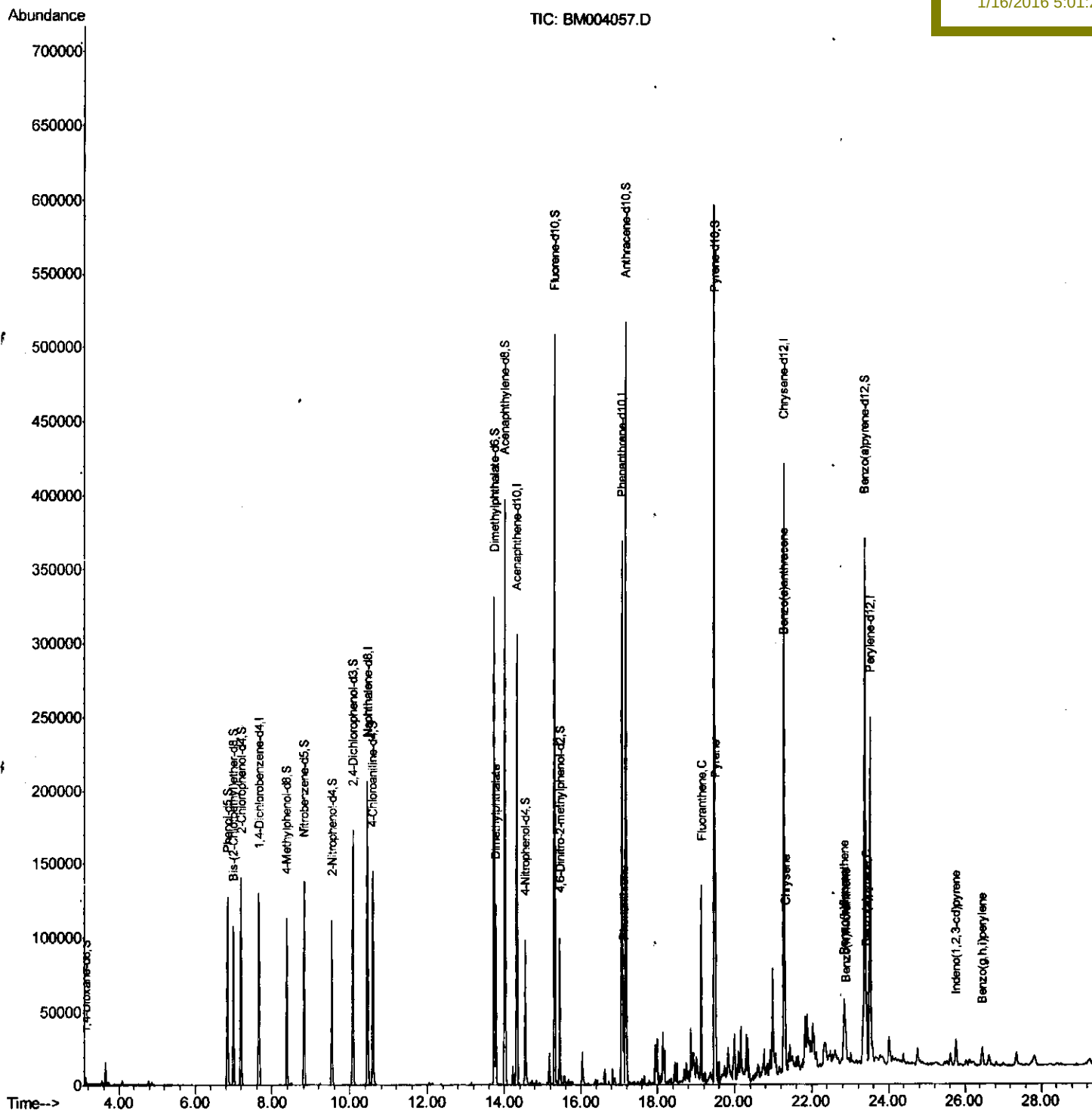
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011616\
Data File : BM004057.D
Acq On : 15 Jan 2016 20:49
Operator : SJ/UM
Sample : H1100-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BC986

Quant Time: Jan 16 03:31:28 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jan 16 03:00:58 2016
Response via : Initial Calibration

Manual Integrations
APPROVED

Sohil
1/16/2016 5:01:28 AM



Quantitation Report (Qedit)

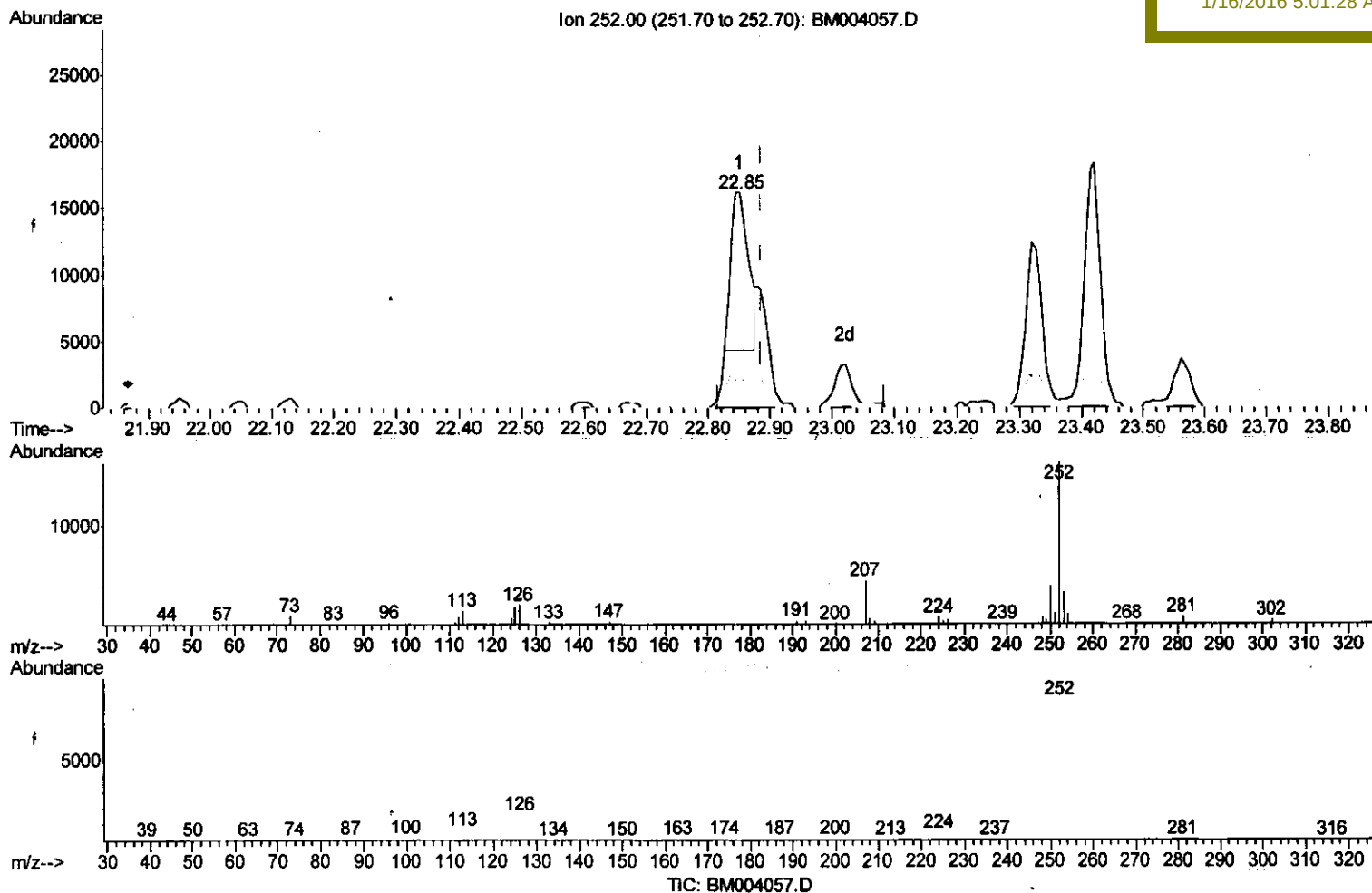
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Sohil
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(89) Benzo(k)fluoranthene

22.850min (-0.035) 2.46ng/ul

response 23179

Ion	Exp%	Act%
252.00	100	100
253.00	21.60	21.62
125.00	10.20	12.45#
0.00	0.00	0.00

Quantitation Report (Qedit)

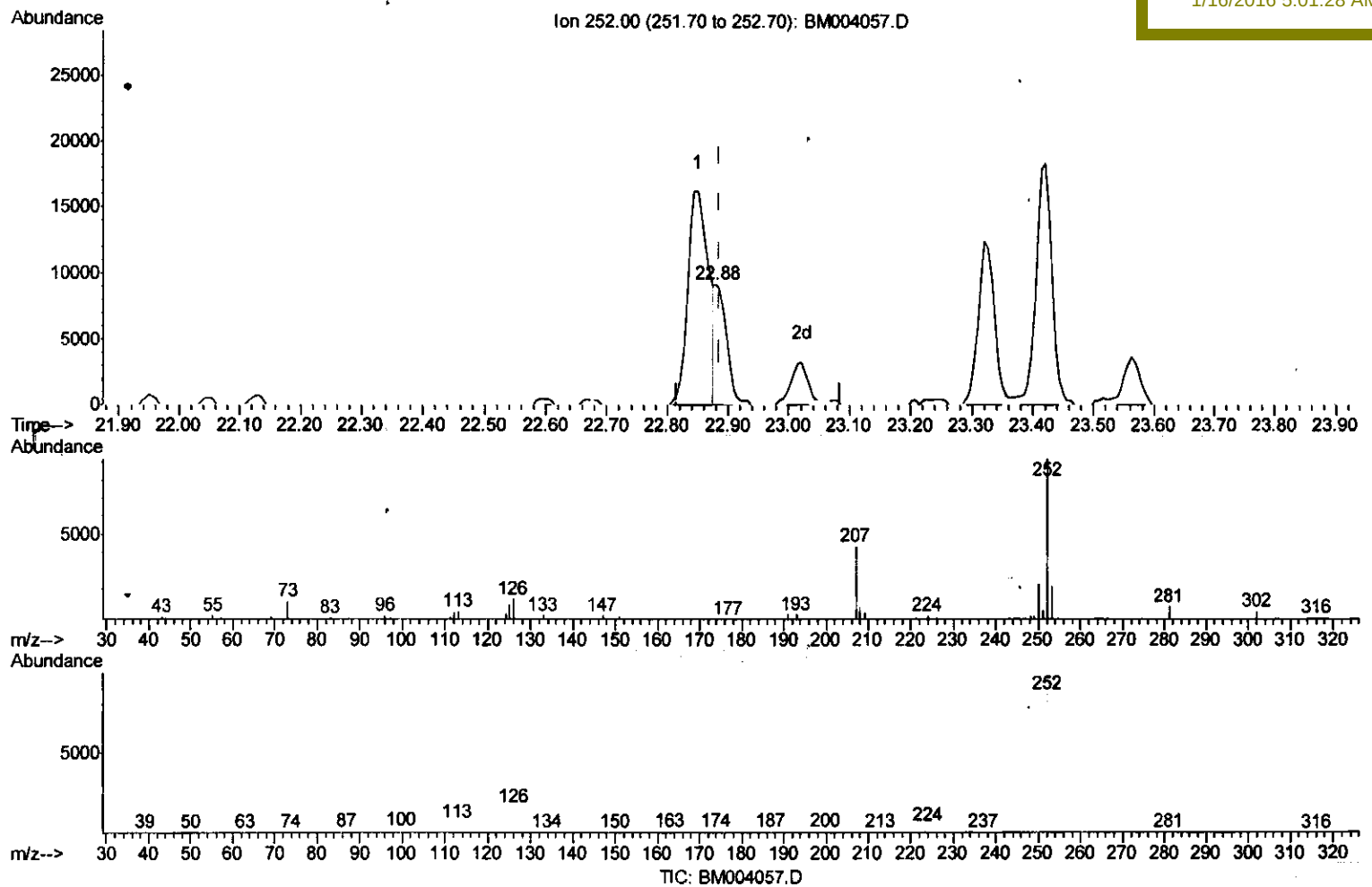
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(89) Benzo(k)fluoranthene

22.880min (-0.006) 1.36ng/ul m

response 12827

Ion Exp% Act%

252.00 100 100

253.00 21.60 23.19

125.00 10.20 12.27#

0.00 0.00 0.00

S.J
 1/16/16

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Internal Standards	R.T.	Q Ion	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	36105	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	175589	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	101051	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	220694	20.00	ng/ul	0.00
f 78) Chrysene-d12	21.28	240	183559	20.00	ng/ul	0.00
86) Perylene-d12	23.51	264	163000	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	3292	4.08	ng/uL	0.00
5) Phenol-d5	6.85	99	79291	26.45	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.00	67	46968	27.72	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	67953	27.58	ng/ul	0.00
13) 4-Methylphenol-d8	8.37	113	46742	19.24	ng/ul	0.00
19) Nitrobenzene-d5	8.82	128	36389	27.51	ng/ul	0.00
22) 2-Nitrophenol-d4	9.54	143	39975	27.57	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	66549	25.83	ng/ul	0.00
29) 4-Chloroaniline-d4	10.59	131	87263	25.66	ng/ul	0.00
44) Dimethylphthalate-d6	13.73	166	229391	28.91	ng/ul	0.00
47) Acenaphthylene-d8	14.01	160	281829	29.16	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	30339	21.29	ng/ul	0.00
58) Fluorene-d10	15.32	176	200204	29.75	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	21393	17.19	ng/ul	0.00
71) Anthracene-d10	17.17	188	286797	28.10	ng/ul	0.00
79) Pyrene-d10	19.47	212	302836	31.86	ng/ul	0.00
f 90) Benzo(a)pyrene-d12	23.37	264	242846	29.83	ng/ul	0.00

Target Compounds

						Qvalue
45) Dimethylphthalate	13.78	163	82247	10.17	ng/ul	100
70) Phenanthrene	17.12	178	44069	3.55	ng/ul	99
77) Fluoranthene	19.14	202	73733	5.72	ng/ul	99
80) Pyrene	19.50	202	102383	8.22	ng/ul	99
83) Benzo(a)anthracene	21.26	228	36127	3.31	ng/ul	96
85) Chrysene	21.32	228	45859	4.42	ng/ul	98
88) Benzo(b)fluoranthene	22.85	252	37939	3.83	ng/ul	98
89) Benzo(k)fluoranthene	22.88	252	12827m	1.36	ng/ul	98
91) Benzo(a)pyrene	23.42	252	33305	3.54	ng/ul	98
92) Indeno(1,2,3-cd)pyrene	25.77	276	18607	1.79	ng/ul	97
94) Benzo(g,h,i)perylene	26.46	276	16522	1.89	ng/ul	99

3.5
 1/16/16

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