

Quantitation Report (Qedit)

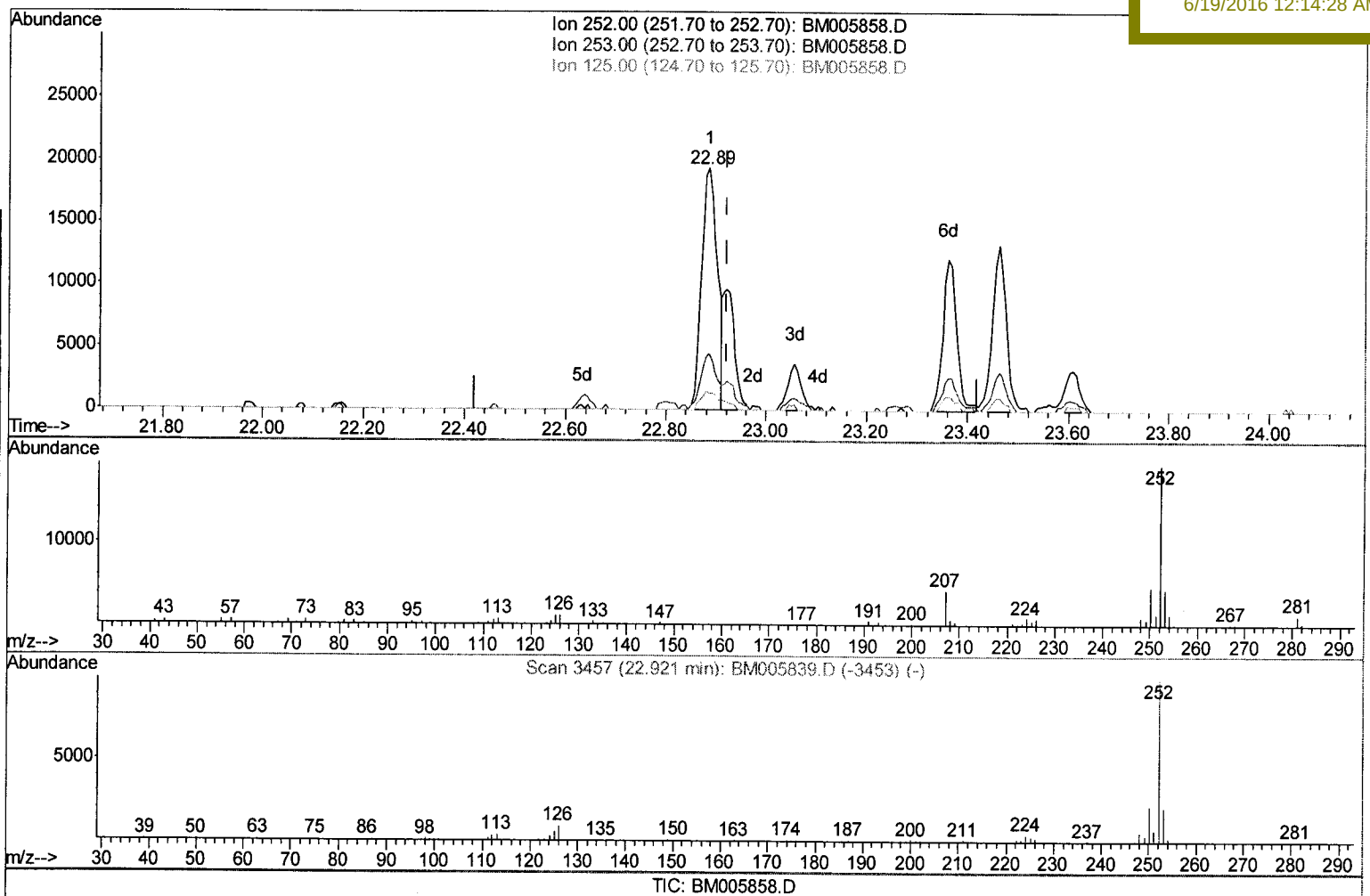
Data Path : Z:\HPCHEM1\BNA_M\Data\BM061816\
 Data File : BM005858.D
 Acq On : 17 Jun 2016 12:58
 Operator : UM/SJ
 Sample : H3535-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9XW9

Quant Time: Jun 17 14:40:18 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061716.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 17 10:48:37 2016
 Response via : Initial Calibration

Manual Integrations
 APPROVED

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 6/19/2016 12:14:28 AM



(36) Benzo(k)fluoranthene
 22.885min (-0.035) 4.02ng/ul
 response 41784

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	23.33
125.00	6.00	6.82
0.00	0.00	0.00

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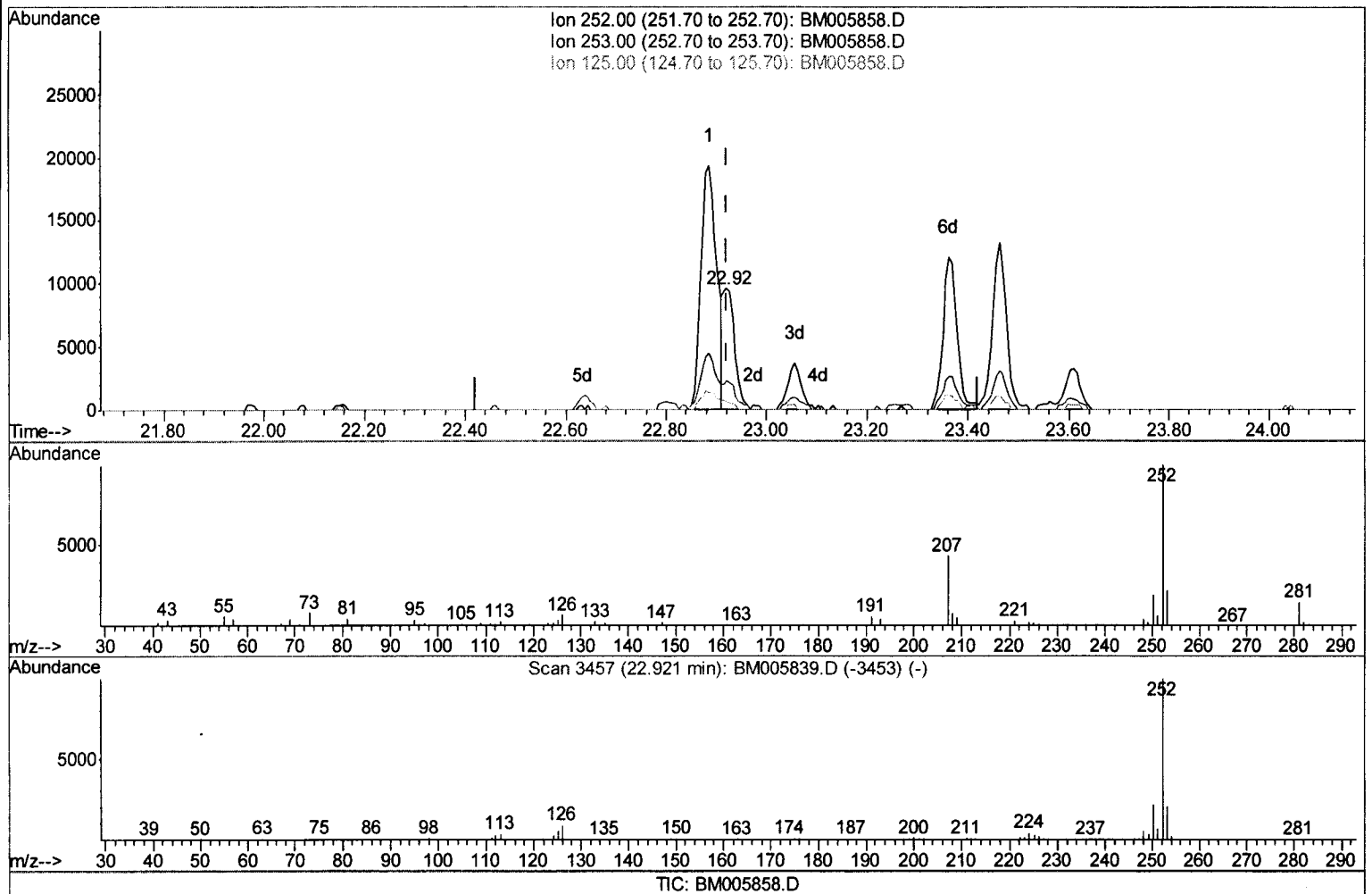
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(36) Benzo(k)fluoranthene

22.921min (-0.000) 1.49ng/ul m

response 15499

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	24.29
125.00	6.00	6.38
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	7.73	152	23779	20.00	ng/ul	0.00
7) Naphthalene-d8	10.52	136	113848	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.37	164	74346	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.12	188	173979	20.00	ng/ul	0.00
29) Chrysene-d12	21.30	240	200015	20.00	ng/ul	0.00
34) Perylene-d12	23.56	264	172727	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.17	96	254	1.38	ng/uL	0.00
3) Phenol-d5	6.92	99	20319	10.21	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.06	67	13102	12.70	ng/ul	0.00
5) 2-Chlorophenol-d4	7.27	132	20586	12.40	ng/ul	0.00
6) 4-Methylphenol-d8	8.45	113	13565	7.62	ng/ul	0.00
8) Nitrobenzene-d5	8.89	128	10570	13.15	ng/ul	0.00
9) 2-Nitrophenol-d4	9.62	143	12368	12.60	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.16	165	22129	11.41	ng/ul	0.00
12) 4-Chloroaniline-d4	10.67	131	5381	4.10	ng/ul	0.00
16) Dimethylphthalate-d6	13.78	166	83834	11.92	ng/ul	0.00
17) Acenaphthylene-d8	14.07	160	106569	13.15	ng/ul	0.00
20) 4-Nitrophenol-d4	14.64	143	1782	2.10	ng/ul	0.03
21) Fluorene-d10	15.37	176	76982	12.84	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.52	200	5558	4.73	ng/ul	0.00
26) Anthracene-d10	17.22	188	118745	13.07	ng/ul	0.00
30) Pyrene-d10	19.52	212	141091	14.88	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.41	264	111896	12.96	ng/ul	0.00

Target Compounds

						Qvalue
13) 2-Methylnaphthalene	12.19	142	4845	1.04	ng/ul	96
25) Phenanthrene	17.16	178	33621	3.25	ng/ul	98
28) Fluoranthene	19.18	202	61829	4.75	ng/ul	98
31) Pyrene	19.54	202	51939	4.39	ng/ul	99
32) Benzo(a)anthracene	21.29	228	27441	2.20	ng/ul	96
33) Chrysene	21.34	228	33919	2.82	ng/ul	97
35) Benzo(b)fluoranthene	22.89	252	41784	3.89	ng/ul	96
36) Benzo(k)fluoranthene	22.92	252	15499m	1.49	ng/ul	
38) Benzo(a)pyrene	23.46	252	25418	2.41	ng/ul#	98
39) Indeno(1,2,3-cd)pyrene	25.84	276	18961	1.47	ng/ul	93
41) Benzo(g,h,i)perylene	26.54	276	19492	1.77	ng/ul	100

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

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