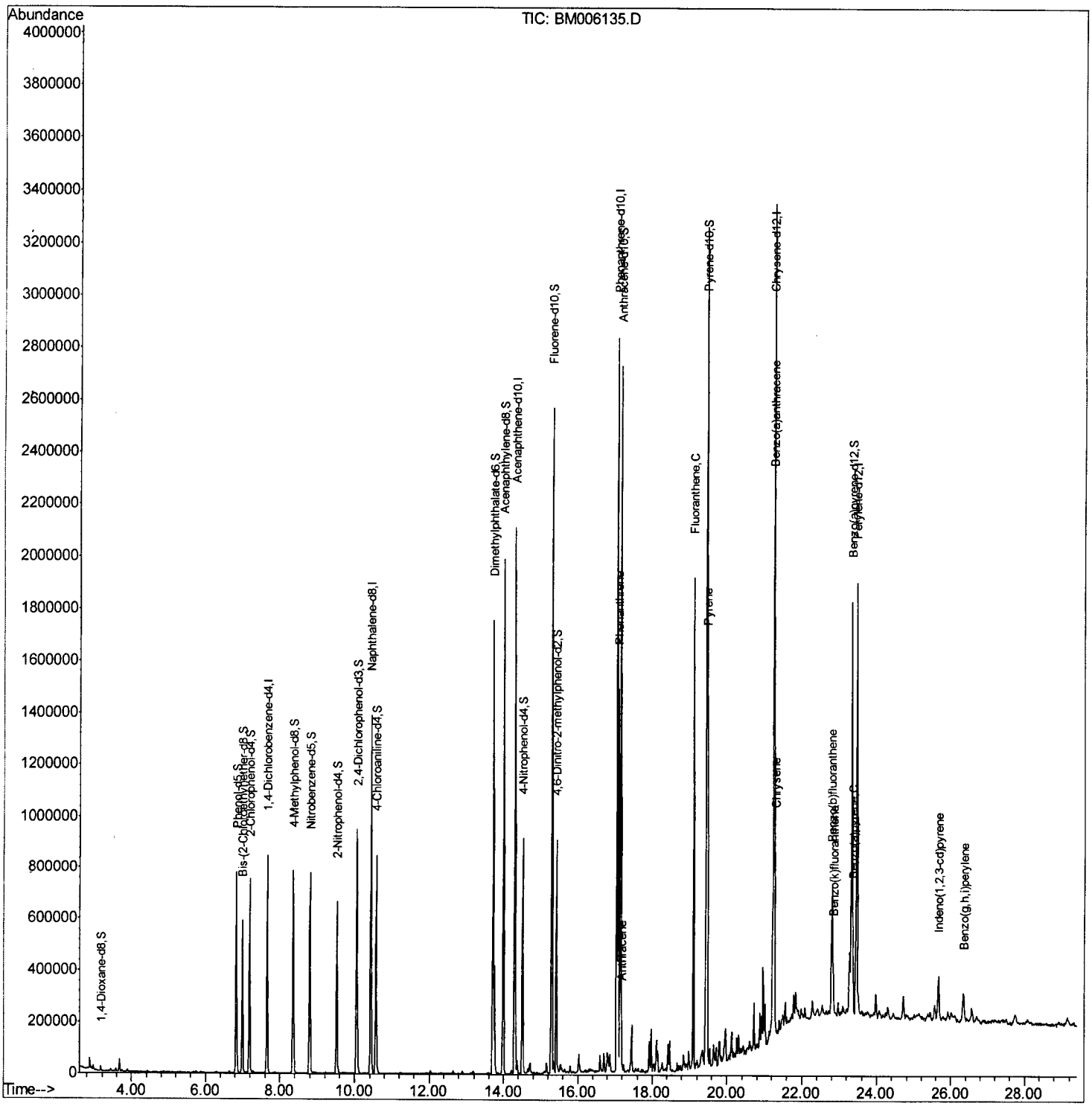
Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062916\
Data File : BM006135.D
Acq On : 29 Jun 2016 21:03
Operator : UM/SJ
Sample : H3779-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Quant Time: Jun 30 03:52:44 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

Instrument :
BNA_M
ClientSampleId :
D9YD8

Manual Integratio
APPROVE
sohil
6/30/2016 7:44:11 PM



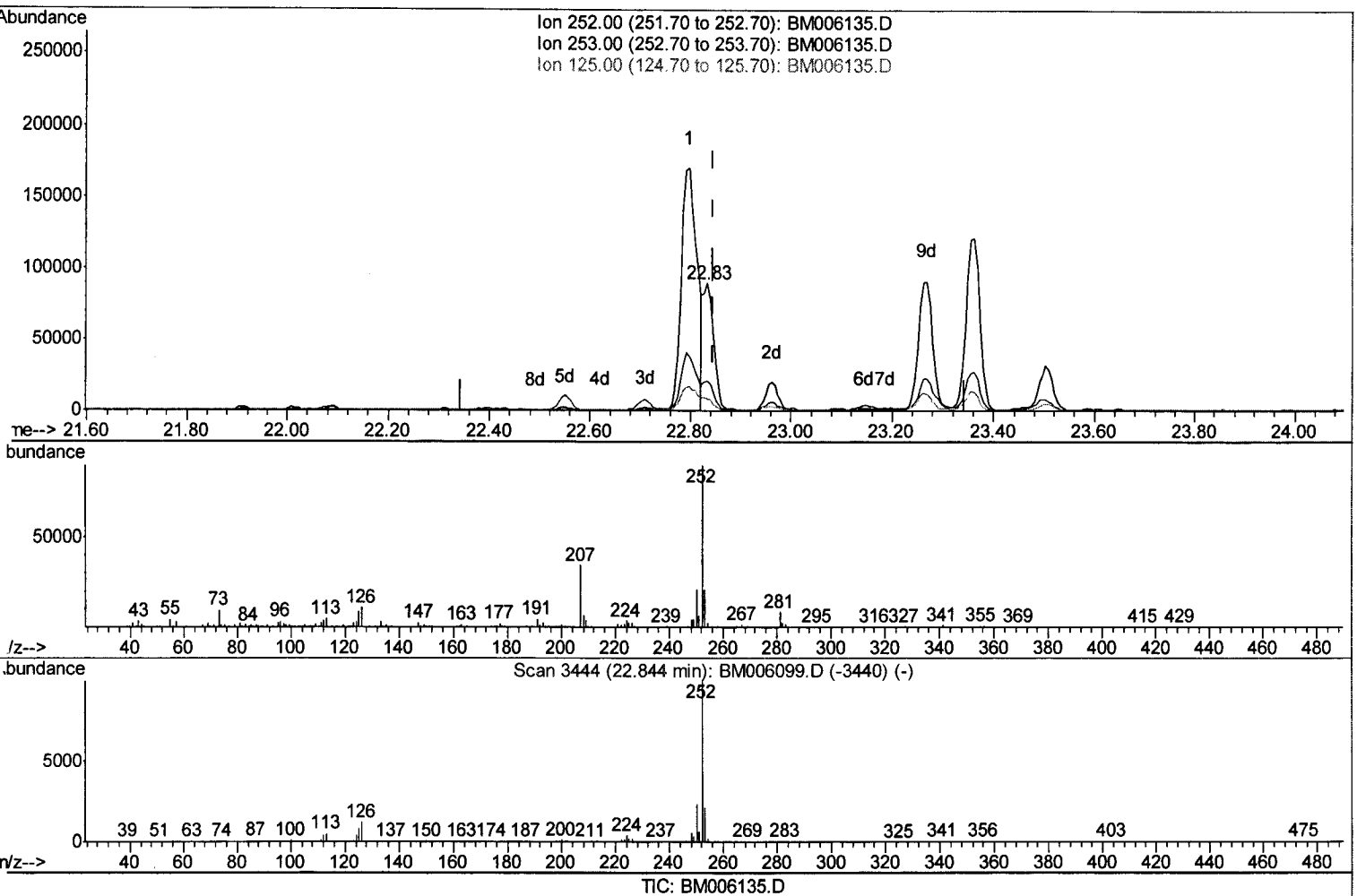
Quantitation Report (Qedit)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062916\
Data File : BM006135.D
Acq On : 29 Jun 2016 21:03
Operator : UM/SJ
Sample : H3779-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Quant Time: Jun 30 03:30:45 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

Instrument :
BNA_M
ClientSampleId :
D9YD8

Manual Integratio
APPROVE
sohil
6/30/2016 7:44:11 PM



TIC: BM006135.D

(36) Benzo(k)fluoranthene

22.832min (-0.012) 2.13ng/ul m V.M. 07/02/16

response 126121

Ion	Exp%	Act%
252.00	100	100
253.00	22.00	23.30
125.00	10.10	9.78
0.00	0.00	0.00

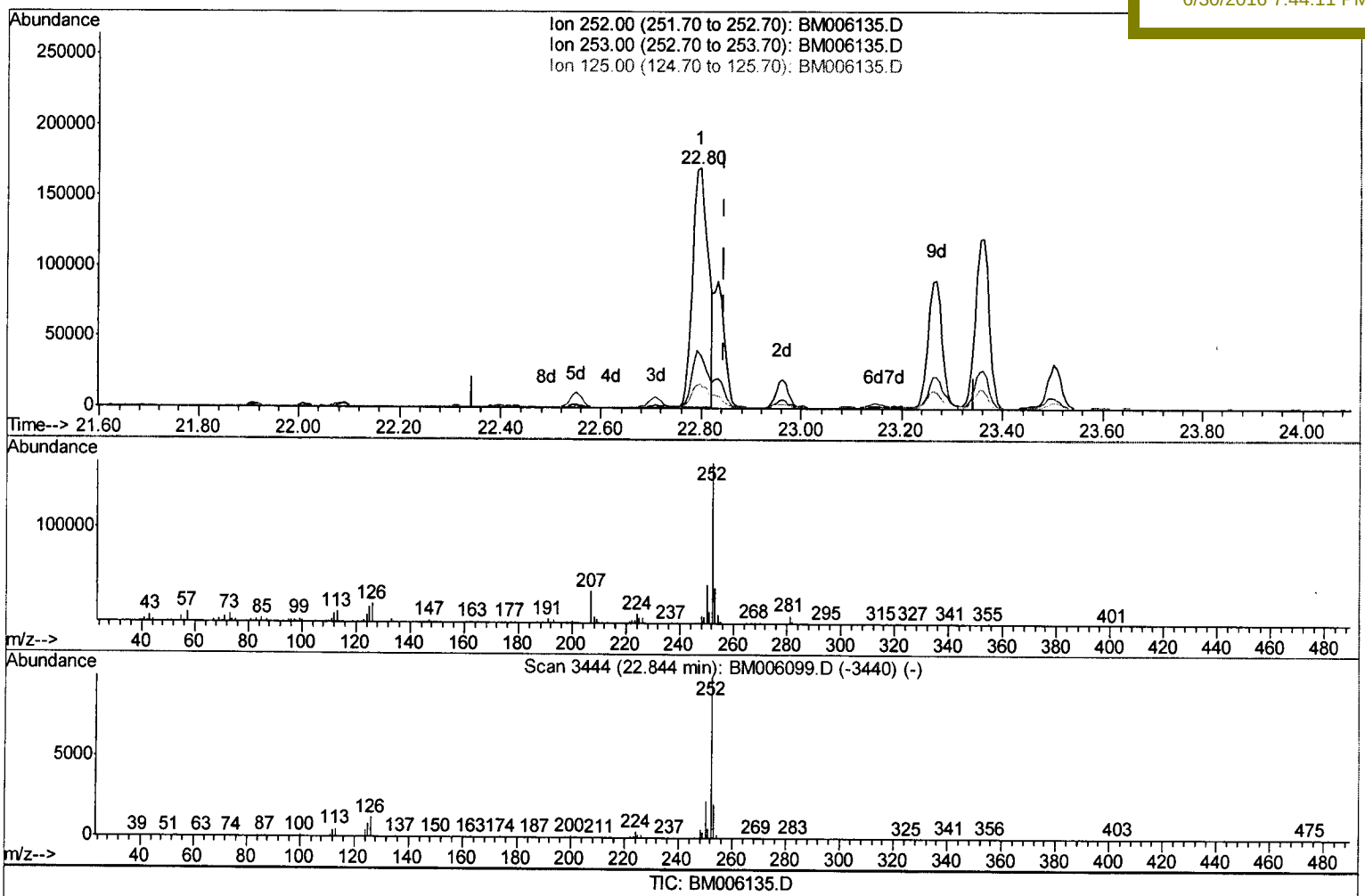
Quantitation Report (Qedit)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062916\
Data File : BM006135.D
Acq On : 29 Jun 2016 21:03
Operator : UM/SJ
Sample : H3779-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Quant Time: Jun 30 03:30:45 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

Instrument :
BNA_M
ClientSampleId :
D9YD8

Manual Integratio
APPROVE
sohil
6/30/2016 7:44:11 PM



TIC: BM006135.D

(36) Benzo(k)fluoranthene

22.797min (-0.047) 6.28ng/ul

response 372147

Ion	Exp%	Act%
252.00	100	100
253.00	22.00	22.42
125.00	10.10	10.01
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062916\
Data File : BM006135.D
Acq On : 29 Jun 2016 21:03
Operator : UM/SJ
Sample : H3779-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Quant Time: Jun 30 03:52:44 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

Instrument : BNA_M
ClientSampleId : D9YD8

Manual Integratio
APPROVE
sohil
6/30/2016 7:44:11 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	221216	20.00	ng/ul	0.00
7) Naphthalene-d8	10.43	136	1154664	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.30	164	735366	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.05	188	1670666	20.00	ng/ul	0.00
29) Chrysene-d12	21.24	240	1445539	20.00	ng/ul	0.00
34) Perylene-d12	23.46	264	1216593	20.00	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.16	96	8328	1.75	ng/uL	0.00
3) Phenol-d5	6.82	99	413953	22.91	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.99	67	266988	25.06	ng/ul	0.00
5) 2-Chlorophenol-d4	7.18	132	322690	23.26	ng/ul	0.00
6) 4-Methylphenol-d8	8.35	113	285531	19.63	ng/ul	-0.01
8) Nitrobenzene-d5	8.80	128	171592	22.06	ng/ul	0.00
9) 2-Nitrophenol-d4	9.52	143	195968	22.02	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.05	165	340731	20.98	ng/ul	0.00
12) 4-Chloroaniline-d4	10.57	131	431377	33.60	ng/ul	0.00
16) Dimethylphthalate-d6	13.71	166	1184173	22.46	ng/ul	0.00
17) Acenaphthylene-d8	13.99	160	1443624	22.57	ng/ul	0.00
20) 4-Nitrophenol-d4	14.50	143	200976	20.09	ng/ul	0.00
21) Fluorene-d10	15.30	176	1027731	22.85	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.42	200	164203	27.75	ng/ul	0.00
26) Anthracene-d10	17.14	188	1507251	22.49	ng/ul	0.00
30) Pyrene-d10	19.44	212	1643451	28.23	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.31	264	1128709	23.31	ng/ul	0.00
Target Compounds						Qvalue
25) Phenanthrene	17.09	178	877432	11.10	ng/ul	100
27) Anthracene	17.18	178	96971	1.19	ng/ul	98
28) Fluoranthene	19.11	202	1154891	11.93	ng/ul	99
31) Pyrene	19.47	202	887037	11.63	ng/ul	98
32) Benzo(a)anthracene	21.23	228	339106	4.51	ng/ul	98
33) Chrysene	21.28	228	367025	5.30	ng/ul	98
35) Benzo(b)fluoranthene	22.80	252	372147	5.89	ng/ul	98
36) Benzo(k)fluoranthene	22.83	252	126121m	2.13	ng/ul	
38) Benzo(a)pyrene	23.36	252	226336	3.79	ng/ul	99
39) Indeno(1,2,3-cd)pyrene	25.67	276	159595	2.57	ng/ul#	92
41) Benzo(g,h,i)perylene	26.34	276	131357	2.53	ng/ul	99

U.M. 07/02/16

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