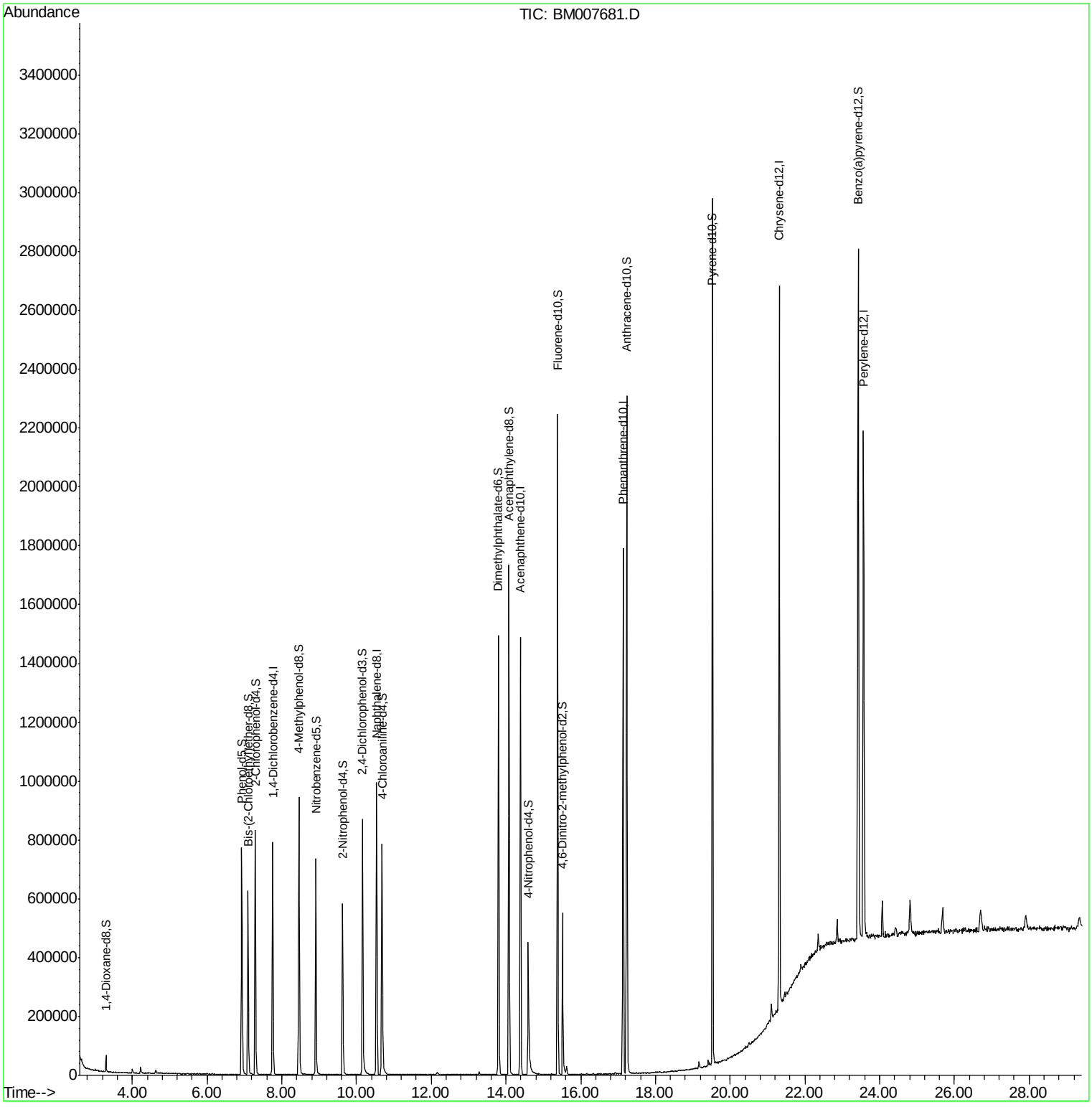
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101916\
Data File : BM007681.D
Acq On : 19 Oct 2016 20:39
Operator : UM/SJ
Sample : MDL-BLK-S-ML
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
MDL-BLK-S-ML

Quant Time: Oct 20 03:12:38 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM101916-MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Oct 20 03:12:33 2016
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101916\
Data File : BM007681.D
Acq On : 19 Oct 2016 20:39
Operator : UM/SJ
Sample : MDL-BLK-S-ML
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : MDL-BLK-S-ML

Quant Time: Oct 20 03:12:38 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM101916-MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Oct 20 03:12:33 2016
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	209240	20.00	ng/ul	0.00
7) Naphthalene-d8	10.53	136	803526	20.00	ng/ul	0.00
13) Acenaphthene-d10	14.39	164	514584	20.00	ng/ul	0.00
18) Phenanthrene-d10	17.13	188	1015545	20.00	ng/ul	0.00
23) Chrysene-d12	21.32	240	1286065	20.00	ng/ul	0.00
25) Perylene-d12	23.57	264	1323154	20.00	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.29	96	27877	5.68	ng/uL	0.00
3) Phenol-d5	6.93	99	450568	27.79	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.09	67	274398	28.56	ng/ul	0.00
5) 2-Chlorophenol-d4	7.29	132	363058	29.17	ng/ul	0.00
6) 4-Methylphenol-d8	8.46	113	352710	27.68	ng/ul	0.00
8) Nitrobenzene-d5	8.91	128	164488	28.98	ng/ul	0.00
9) 2-Nitrophenol-d4	9.63	143	181751	29.34	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.16	165	319770	26.23	ng/ul	0.00
11) 4-Chloroaniline-d4	10.68	131	442179	31.96	ng/ul	0.00
14) Dimethylphthalate-d6	13.80	166	1020481	28.43	ng/ul	0.00
15) Acenaphthylene-d8	14.08	160	1215104	28.87	ng/ul	0.00
16) 4-Nitrophenol-d4	14.60	143	127000	21.51	ng/ul	0.00
17) Fluorene-d10	15.38	176	876768	28.30	ng/ul	0.00
19) 4,6-Dinitro-2-methylphenol	15.52	200	137560	22.35	ng/ul	0.00
20) Anthracene-d10	17.23	188	1320655	28.60	ng/ul	0.00
24) Pyrene-d10	19.53	212	1543639	29.65	ng/ul	0.00
26) Benzo(a)pyrene-d12	23.43	264	1663941	28.87	ng/ul	0.00

Target Compounds	Qvalue

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