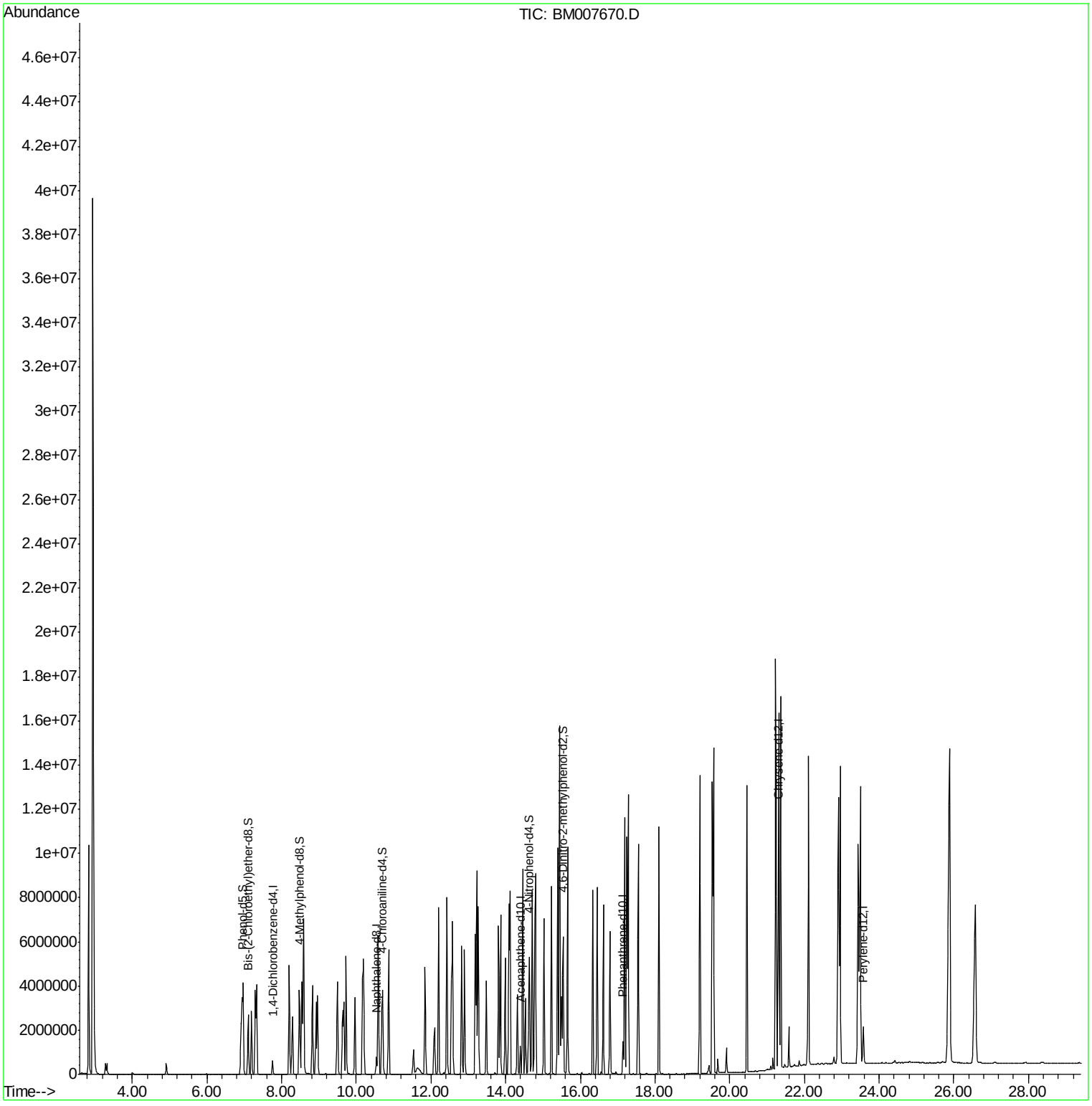
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101916\
Data File : BM007670.D
Acq On : 19 Oct 2016 14:07
Operator : UM/SJ
Sample : SSTD16039
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SSTD16039

Quant Time: Oct 19 14:52:30 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM101916-MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 19 13:45:45 2016
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101916\
Data File : BM007670.D
Acq On : 19 Oct 2016 14:07
Operator : UM/SJ
Sample : SSTD16039
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTD16039

Quant Time: Oct 19 14:52:30 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM101916-MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 19 13:45:45 2016
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	162072	20.00	ng/ul	0.00
7) Naphthalene-d8	10.54	136	647907	20.00	ng/ul	0.00
13) Acenaphthene-d10	14.39	164	431198	20.00	ng/ul	0.00
18) Phenanthrene-d10	17.14	188	891889	20.00	ng/ul	0.00
23) Chrysene-d12	21.33	240	1121806	20.00	ng/ul	0.00
25) Perylene-d12	23.58	264	1183370	20.00	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
3) Phenol-d5	6.94	99	2149471	140.40	ng/ul	0.01
4) Bis-(2-Chloroethyl)ether-d	7.10	67	1194046	148.42	ng/ul	0.01
5) 2-Chlorophenol-d4	0.00	132	0d	0.00	ng/ul	
6) 4-Methylphenol-d8	8.47	113	1691108	126.66	ng/ul	0.01
8) Nitrobenzene-d5	0.00	128	0d	0.00	ng/ul	
9) 2-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
10) 2,4-Dichlorophenol-d3	0.00	165	0d	0.00	ng/ul	
11) 4-Chloroaniline-d4	10.69	131	1584136	114.84	ng/ul	0.00
14) Dimethylphthalate-d6	0.00	166	0d	0.00	ng/ul	
15) Acenaphthylene-d8	0.00	160	0d	0.00	ng/ul	
16) 4-Nitrophenol-d4	14.63	143	902542	128.52	ng/ul	0.02
17) Fluorene-d10	0.00	176	0d	0.00	ng/ul	
19) 4,6-Dinitro-2-methylphenol	15.53	200	1000828	178.46	ng/ul	0.02
20) Anthracene-d10	0.00	188	0d	0.00	ng/ul	
24) Pyrene-d10	0.00	212	0d	0.00	ng/ul	
26) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul	

Target Compounds	Qvalue

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