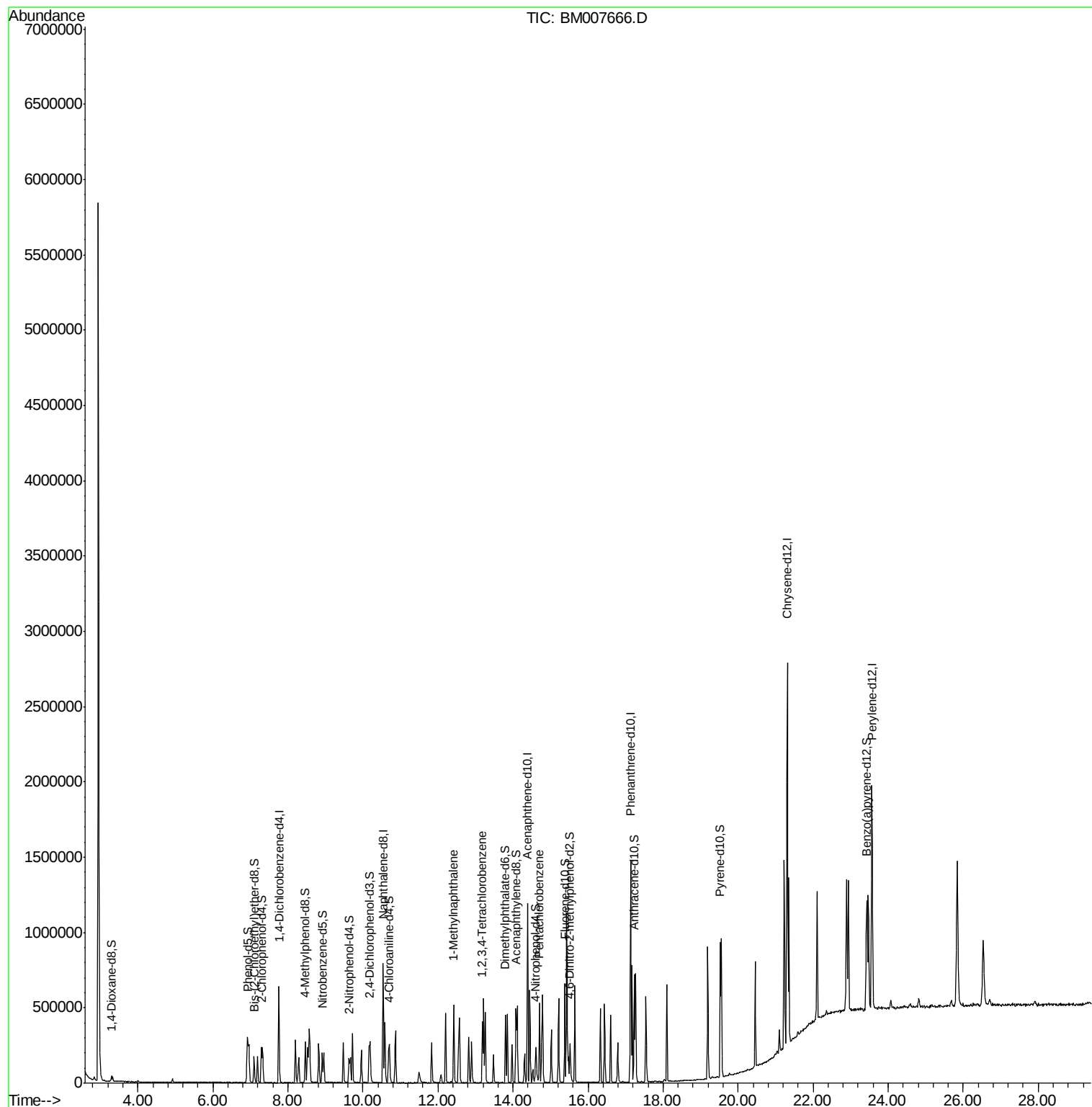


Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101916\
Data File : BM007666.D
Acq On : 19 Oct 2016 11:26
Operator : UM/SJ
Sample : SSTD01035
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SSTD01035

Quant Time: Oct 19 13:48:13 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



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 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	167736	20.00	ng/ul	0.00
7) Naphthalene-d8	10.54	136	642707	20.00	ng/ul	0.00
13) Acenaphthene-d10	14.39	164	417120	20.00	ng/ul	0.00
18) Phenanthrene-d10	17.13	188	833401	20.00	ng/ul	0.00
23) Chrysene-d12	21.32	240	1106456	20.00	ng/ul	0.00
25) Perylene-d12	23.57	264	1120012	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.29	96	16116	4.91	ng/uL	0.00
3) Phenol-d5	6.93	99	125287	7.91	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.10	67	78400	9.42	ng/ul	0.00
5) 2-Chlorophenol-d4	7.29	132	99610	8.51	ng/ul	0.00
6) 4-Methylphenol-d8	8.46	113	97074	7.02	ng/ul	0.00
8) Nitrobenzene-d5	8.91	128	44124	9.35	ng/ul	0.00
9) 2-Nitrophenol-d4	9.63	143	48314	8.77	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.16	165	97107	8.85	ng/ul	0.00
11) 4-Chloroaniline-d4	10.68	131	107091	7.83	ng/ul	0.00
14) Dimethylphthalate-d6	13.80	166	290549	8.07	ng/ul	0.00
15) Acenaphthylene-d8	14.08	160	336190	8.06	ng/ul	0.00
16) 4-Nitrophenol-d4	14.60	143	39149	5.76	ng/ul	0.00
17) Fluorene-d10	15.38	176	253528	8.15	ng/ul	0.00
19) 4,6-Dinitro-2-methylphenol	15.52	200	42311	8.07	ng/ul	0.00
20) Anthracene-d10	17.23	188	381885	9.81	ng/ul	0.00
24) Pyrene-d10	19.53	212	449927	9.73	ng/ul	0.00
26) Benzo(a)pyrene-d12	23.43	264	496524	9.64	ng/ul	0.00

Target Compounds

						Qvalue
12) 1-Methylnaphthalene	12.42	142	224717	8.92	ng/ul	98
21) 1,2,3,4-Tetrachlorobenzene	13.19	216	125369	12.46	ng/uL	98
22) Pentachlorobenzene	14.71	250	131344	11.82	ng/uL	97

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