

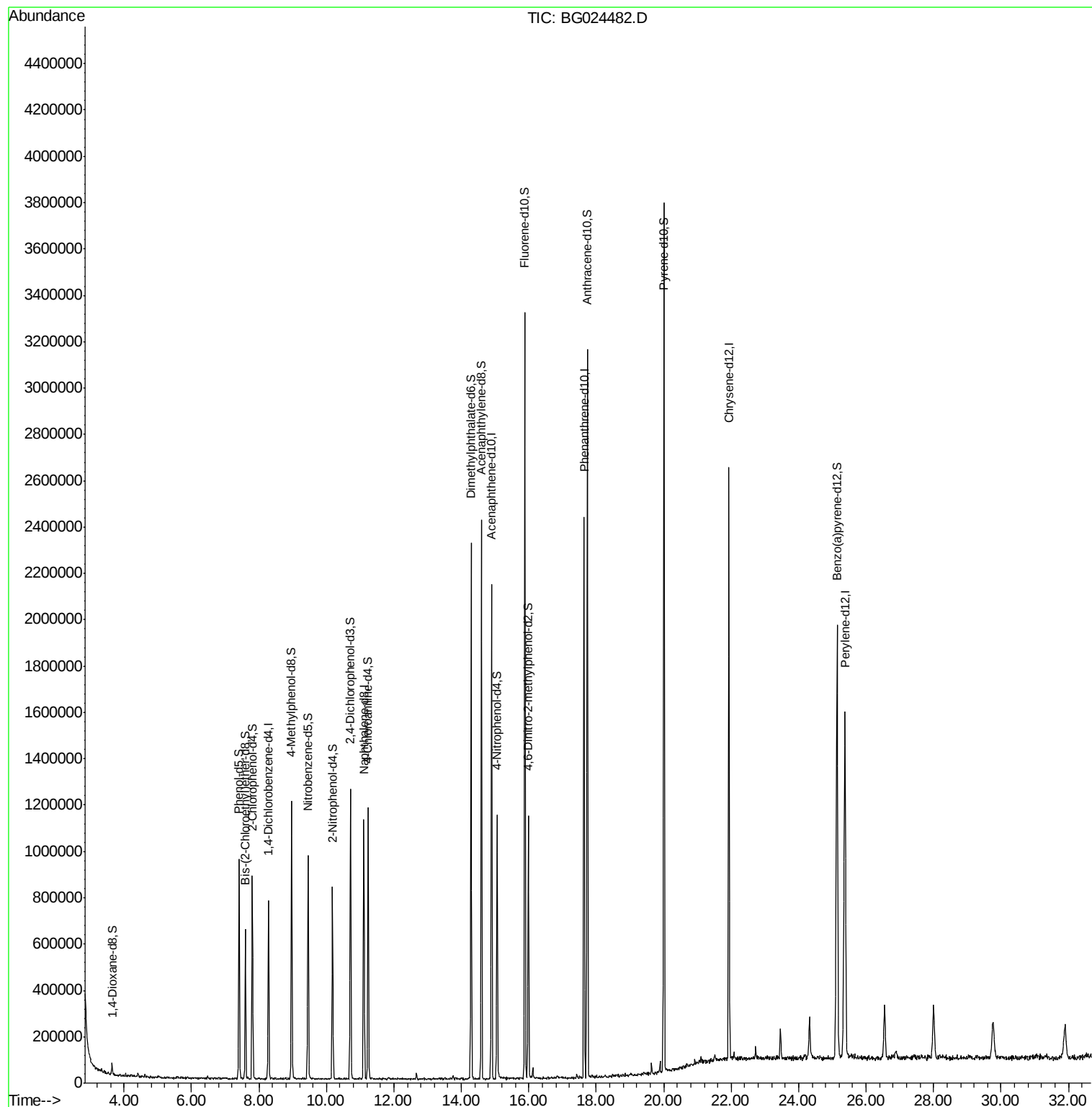
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102016\
Data File : BG024482.D
Acq On : 21 Oct 2016 3:40
Operator : UM/SJ
Sample : MDL-BLK-W
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MDL-BLK-W

Manual Integrations
APPROVED

sohil
10/21/2016 6:54:32 PM

Quant Time: Oct 21 18:12:43 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102016-MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 21 17:27:43 2016
Response via : Initial Calibration



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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.28	152	203485	20.00	ng/ul	0.00
7) Naphthalene-d8	11.11	136	915839	20.00	ng/ul	0.00
13) Acenaphthene-d10	14.89	164	769361	20.00	ng/ul	0.00
18) Phenanthrene-d10	17.63	188	1495454m	20.00	ng/ul	0.00
23) Chrysene-d12	21.93	240	1709601	20.00	ng/ul	0.00
25) Perylene-d12	25.38	264	1795970	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.64	96	23338	6.13	ng/uL	0.00
3) Phenol-d5	7.41	99	494067	26.70	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.59	67	329676	28.82	ng/ul	0.00
5) 2-Chlorophenol-d4	7.80	132	361920	27.79	ng/ul	0.00
6) 4-Methylphenol-d8	8.97	113	435997	28.25	ng/ul	0.00
8) Nitrobenzene-d5	9.46	128	200766	29.31	ng/ul	0.00
9) 2-Nitrophenol-d4	10.18	143	236365	29.64	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.71	165	434995	26.93	ng/ul	0.00
11) 4-Chloroaniline-d4	11.23	131	573230	34.11	ng/ul	0.00
14) Dimethylphthalate-d6	14.29	166	1461273	28.16	ng/ul	0.00
15) Acenaphthylene-d8	14.60	160	1702346	27.57	ng/ul	0.00
16) 4-Nitrophenol-d4	15.06	143	226503	25.54	ng/ul	0.00
17) Fluorene-d10	15.88	176	1340673	27.12	ng/ul	0.00
19) 4,6-Dinitro-2-methylphenol	15.99	200	251290m	25.14	ng/ul	0.00
20) Anthracene-d10	17.73	188	1850607	27.85	ng/ul	0.00
24) Pyrene-d10	20.01	212	2145467	28.89	ng/ul	0.00
26) Benzo(a)pyrene-d12	25.13	264	2239119	28.49	ng/ul	0.00

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