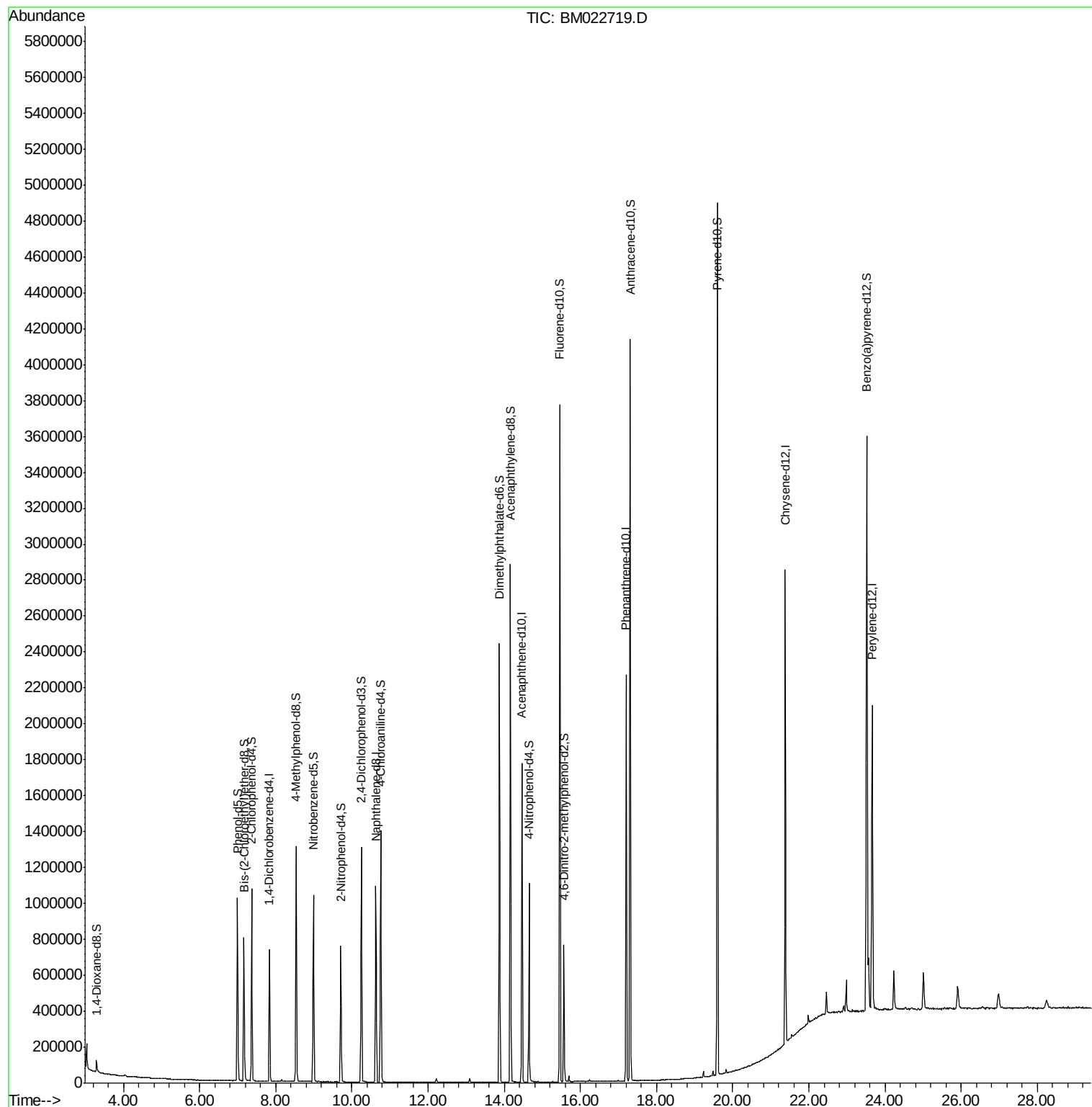


Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM091819\
Data File : BM022719.D
Acq On : 18 Sep 2019 17:26
Operator : HP/JU
Sample : MDL-S-ME-BL
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
MDL-S-ME-BL

Quant Time: Sep 18 18:02:50 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Sep 18 17:31:32 2019
Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM091819\
 Data File : BM022719.D
 Acq On : 18 Sep 2019 17:26
 Operator : HP/JU
 Sample : MDL-S-ME-BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-S-ME-BL

Quant Time: Sep 18 18:02:50 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 18 17:31:32 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	204248	20.00	ng/ul	0.00
18) Naphthalene-d8	10.63	136	912042	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.47	164	583012	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.20	188	1206119	20.00	ng/ul	0.00
78) Chrysene-d12	21.38	240	1120278	20.00	ng/ul	0.00
86) Perylene-d12	23.66	264	1138041	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.29	96	35449	7.23	ng/uL	0.00
5) Phenol-d5	6.99	99	592414	31.44	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.16	67	363037	33.62	ng/ul	0.00
9) 2-Chlorophenol-d4	7.37	132	486433	32.69	ng/ul	0.00
13) 4-Methylphenol-d8	8.53	113	476988	31.14	ng/ul	0.00
19) Nitrobenzene-d5	8.99	128	250571	38.08	ng/ul	0.00
22) 2-Nitrophenol-d4	9.71	143	238381	39.29	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.25	165	470525	30.54	ng/ul	0.00
29) 4-Chloroaniline-d4	10.76	131	741313	36.42	ng/ul	0.00
44) Dimethylphthalate-d6	13.87	166	1622865	34.35	ng/ul	0.00
47) Acenaphthylene-d8	14.16	160	1876582	33.17	ng/ul	0.00
52) 4-Nitrophenol-d4	14.65	143	256474	31.78	ng/ul	0.00
58) Fluorene-d10	15.46	176	1404278	35.50	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.57	200	158000	28.82	ng/ul	0.00
71) Anthracene-d10	17.30	188	2176158	35.22	ng/ul	0.00
79) Pyrene-d10	19.59	212	2397939	38.58	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.52	264	2206776	36.08	ng/ul	0.00

Target Compounds	Qvalue
------------------	--------

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