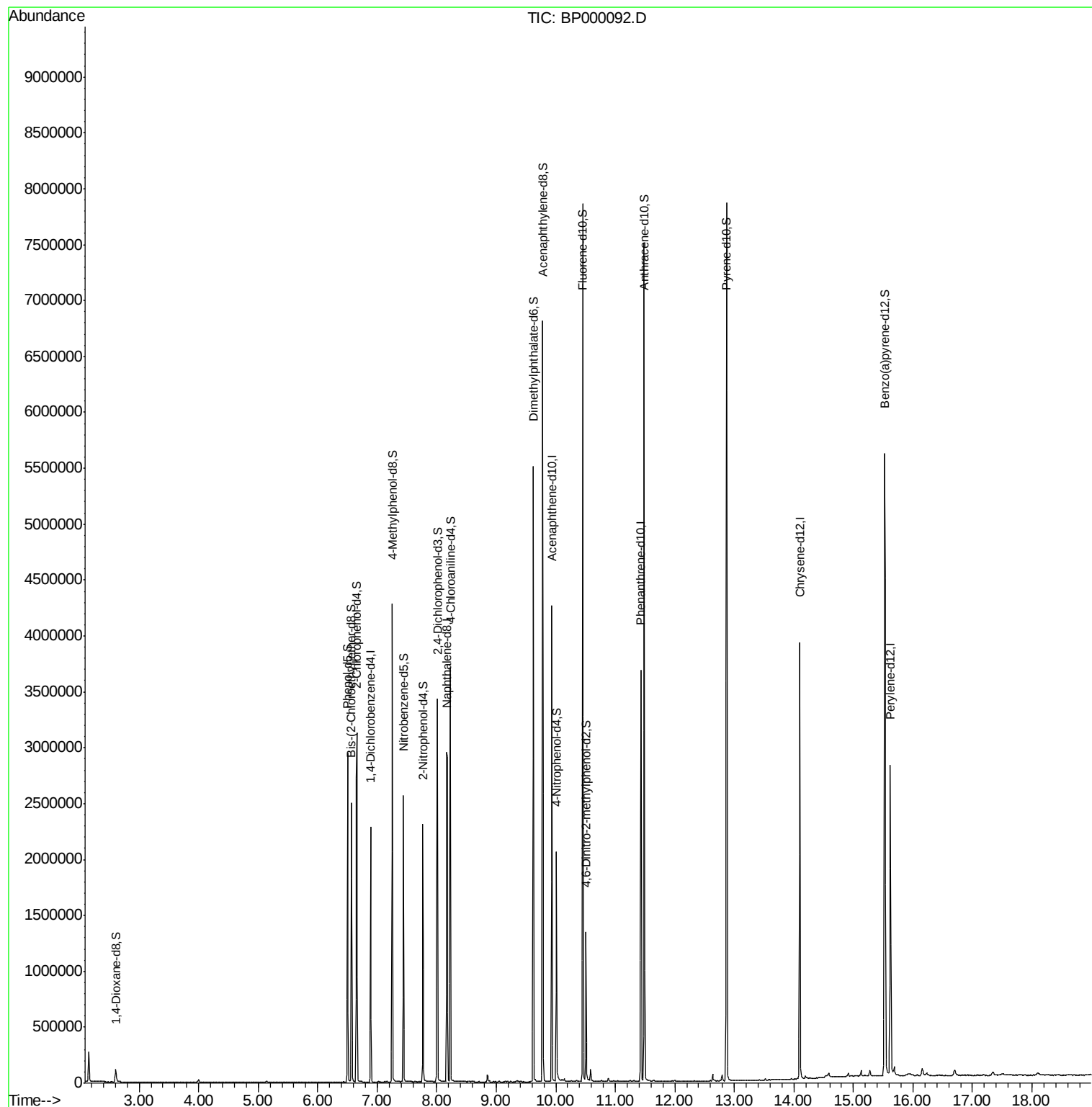


Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090619\
Data File : BP000092.D
Acq On : 06 Sep 2019 16:18
Operator : HP/JU
Sample : MDL-W-BL-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MDL-W-BL-01

Quant Time: Sep 06 17:24:11 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP090519MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Sep 06 01:10:53 2019
Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP090619\
 Data File : BP000092.D
 Acq On : 06 Sep 2019 16:18
 Operator : HP/JU
 Sample : MDL-W-BL-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MDL-W-BL-01

Quant Time: Sep 06 17:24:11 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP090519MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 06 01:10:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	335958	20.00	ng/ul	0.00
18) Naphthalene-d8	8.17	136	1317059	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.93	164	731421	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.43	188	1337156	20.00	ng/ul	0.00
78) Chrysene-d12	14.10	240	1115721	20.00	ng/ul	0.00
86) Perylene-d12	15.62	264	1255849	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.60	96	67695	7.15	ng/uL	0.00
5) Phenol-d5	6.50	99	1026825	35.88	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.57	67	551673	36.87	ng/ul	0.00
9) 2-Chlorophenol-d4	6.66	132	837473	36.19	ng/ul	0.00
13) 4-Methylphenol-d8	7.25	113	763437	35.51	ng/ul	0.00
19) Nitrobenzene-d5	7.45	128	368090	36.81	ng/ul	0.00
22) 2-Nitrophenol-d4	7.77	143	367705	37.77	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	8.01	165	700918	33.92	ng/ul	0.00
29) 4-Chloroaniline-d4	8.23	131	1048498	41.03	ng/ul	0.00
44) Dimethylphthalate-d6	9.62	166	1953842	36.18	ng/ul	0.00
47) Acenaphthylene-d8	9.78	160	2346251	35.15	ng/ul	0.00
52) 4-Nitrophenol-d4	10.01	143	296320	30.71	ng/ul	0.00
58) Fluorene-d10	10.46	176	1662553	36.25	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.51	200	153286	24.94	ng/ul	0.00
71) Anthracene-d10	11.49	188	2414218	37.28	ng/ul	0.00
79) Pyrene-d10	12.87	212	2478186	37.35	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.53	264	2404902	36.94	ng/ul	0.00

Target Compounds Ovalue

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