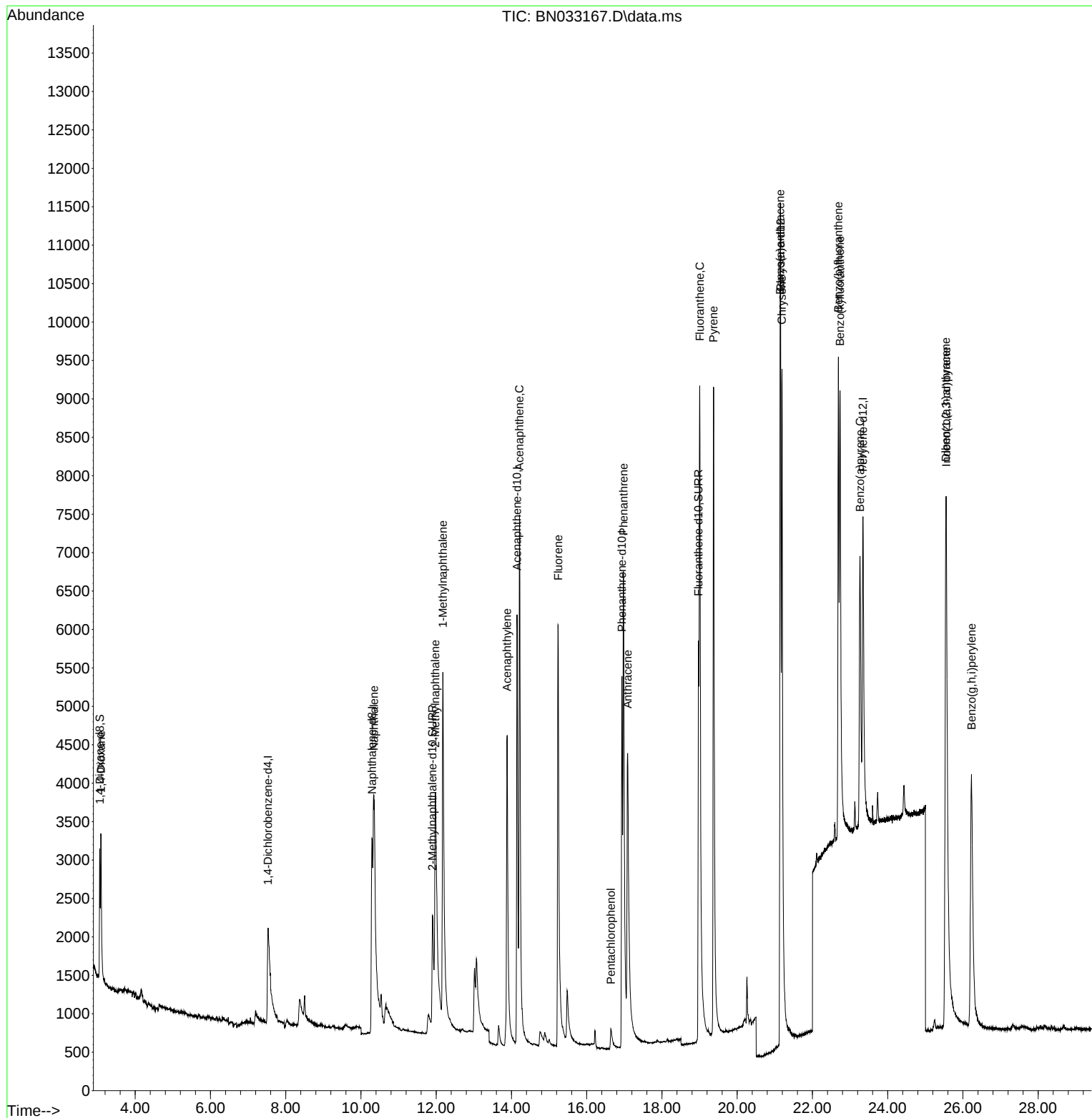


Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080124\
Data File : BN033167.D
Acq On : 01 Aug 2024 13:45
Operator : MA/JU
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SSTD0.4293

Quant Time: Aug 01 16:40:35 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN072524.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 30 14:49:49 2024
Response via : Initial Calibration



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080124\
 Data File : BN033167.D
 Acq On : 01 Aug 2024 13:45
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4293

Quant Time: Aug 01 16:40:35 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN072524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 30 14:49:49 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.529	152	2284	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.295	136	8351	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.148	164	4706	0.400	ng/ul	-0.01
13) Phenanthrene-d10	16.937	188	9550	0.400	ng/ul	0.00
17) Chrysene-d12	21.151	240	7220	0.400	ng/ul	0.00
23) Perylene-d12	23.343	264	7254	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.056	96	1341	0.464	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.901	152	4687	0.344	ng/ul	-0.03
18) Fluoranthene-d10	18.972	212	9608	0.470	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.090	88	1496	0.454	ng/ul#	79
5) Naphthalene	10.344	128	8896	0.415	ng/ul	98
7) 2-Methylnaphthalene	11.978	142	4910	0.338	ng/ul	94
8) 1-Methylnaphthalene	12.176	142	5382	0.343	ng/ul	98
10) Acenaphthylene	13.884	152	7841	0.407	ng/ul	99
11) Acenaphthene	14.213	153	6245	0.431	ng/ul	98
12) Fluorene	15.240	166	6676	0.395	ng/ul	98
14) Pentachlorophenol	16.641	266	466	0.240	ng/ul	96
15) Phenanthrene	16.979	178	9801	0.383	ng/ul	99
16) Anthracene	17.089	178	8211	0.357	ng/ul	99
19) Fluoranthene	19.005	202	11746	0.494	ng/ul	98
20) Pyrene	19.372	202	12140	0.476	ng/ul	99
21) Benzo(a)anthracene	21.140	228	9078	0.369	ng/ul	98
22) Chrysene	21.189	228	12481	0.434	ng/ul	98
24) Benzo(b)fluoranthene	22.688	252	8816	0.416	ng/ul	94
25) Benzo(k)fluoranthene	22.732	252	11401	0.449	ng/ul#	94
26) Benzo(a)pyrene	23.264	252	8077	0.405	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	25.557	276	12016	0.396	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.547	278	9407	0.391	ng/ul	96
29) Benzo(g,h,i)perylene	26.228	276	9914	0.411	ng/ul	97

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