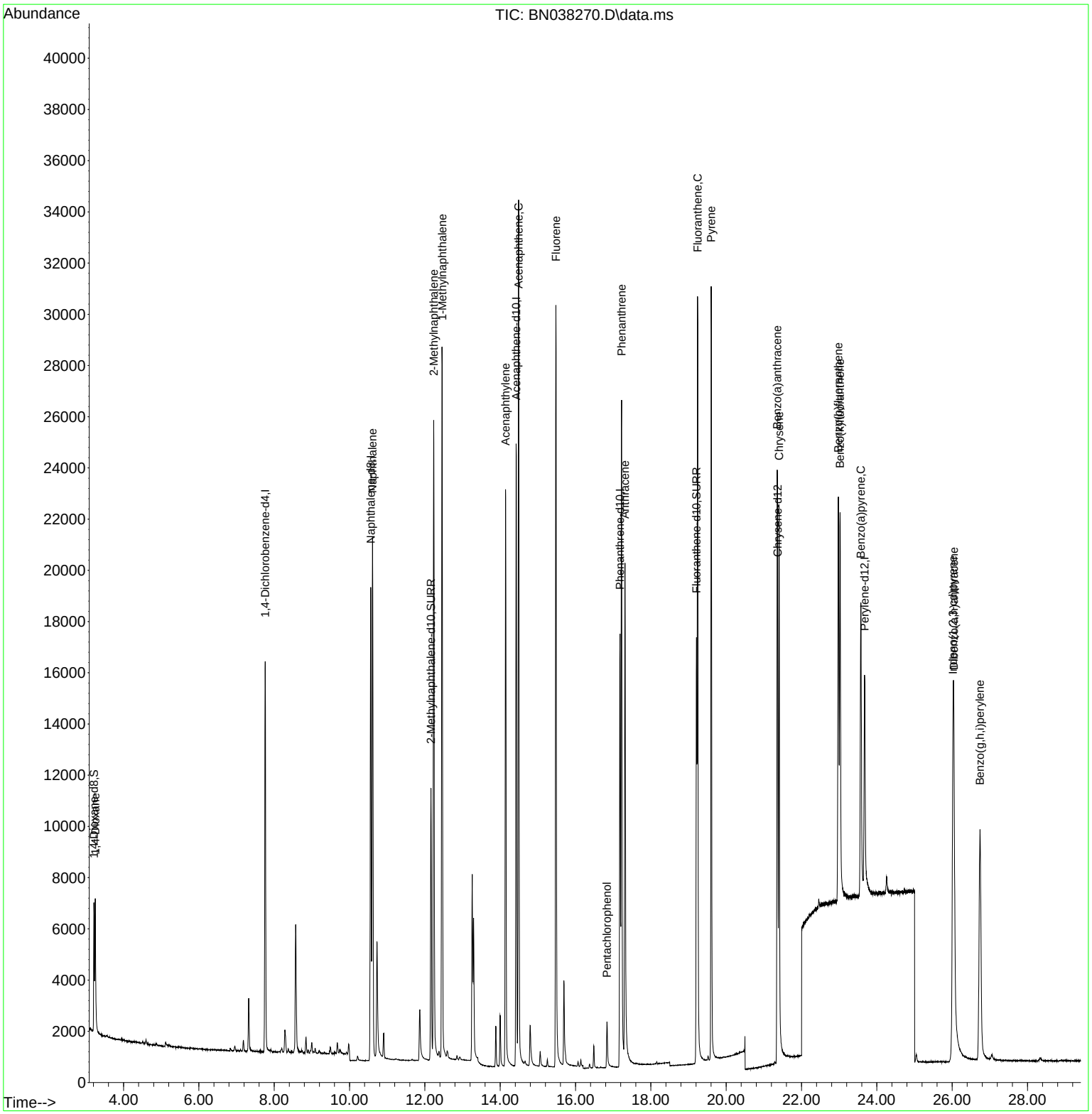
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120525\
Data File : BN038270.D
Acq On : 05 Dec 2025 12:05
Operator : RC/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SSTD0.4277

Quant Time: Dec 05 12:36:14 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN112625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 01 16:36:13 2025
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120525\
Data File : BN038270.D
Acq On : 05 Dec 2025 12:05
Operator : RC/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : SSTD0.4277

Quant Time: Dec 05 12:36:14 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN112625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 01 16:36:13 2025
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.762	152	8084	0.400	ng/ul	0.00
4) Naphthalene-d8	10.562	136	25043	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.424	164	13865	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.180	188	22534	0.400	ng/ul	0.00
17) Chrysene-d12	21.369	240	12432	0.400	ng/ul	0.00
23) Perylene-d12	23.675	264	12699	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.214	96	3383	0.399	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.163	152	14596	0.409	ng/ul	-0.01
18) Fluoranthene-d10	19.212	212	19218	0.398	ng/ul	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.248	88	3526	0.386	ng/ul	99
5) Naphthalene	10.612	128	28289	0.416	ng/ul	100
7) 2-Methylnaphthalene	12.234	142	19051	0.419	ng/ul	100
8) 1-Methylnaphthalene	12.460	142	20436	0.424	ng/ul	99
10) Acenaphthylene	14.142	152	27352	0.420	ng/ul	99
11) Acenaphthene	14.489	153	20088	0.415	ng/ul	99
12) Fluorene	15.479	166	20925	0.403	ng/ul	99
14) Pentachlorophenol	16.834	266	1632	0.452	ng/ul	99
15) Phenanthrene	17.223	178	29796	0.422	ng/ul	100
16) Anthracene	17.316	178	25202	0.420	ng/ul	99
19) Fluoranthene	19.240	202	27513	0.414	ng/ul	98
20) Pyrene	19.603	202	27124	0.412	ng/ul	97
21) Benzo(a)anthracene	21.354	228	18419	0.443	ng/ul	100
22) Chrysene	21.404	228	21603	0.419	ng/ul	99
24) Benzo(b)fluoranthene	22.979	252	21286	0.428	ng/ul	97
25) Benzo(k)fluoranthene	23.023	252	21515	0.411	ng/ul	96
26) Benzo(a)pyrene	23.576	252	19672	0.438	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.022	276	25082	0.469	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.045	278	18437	0.492	ng/ul	97
29) Benzo(g,h,i)perylene	26.739	276	21679	0.431	ng/ul	100

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