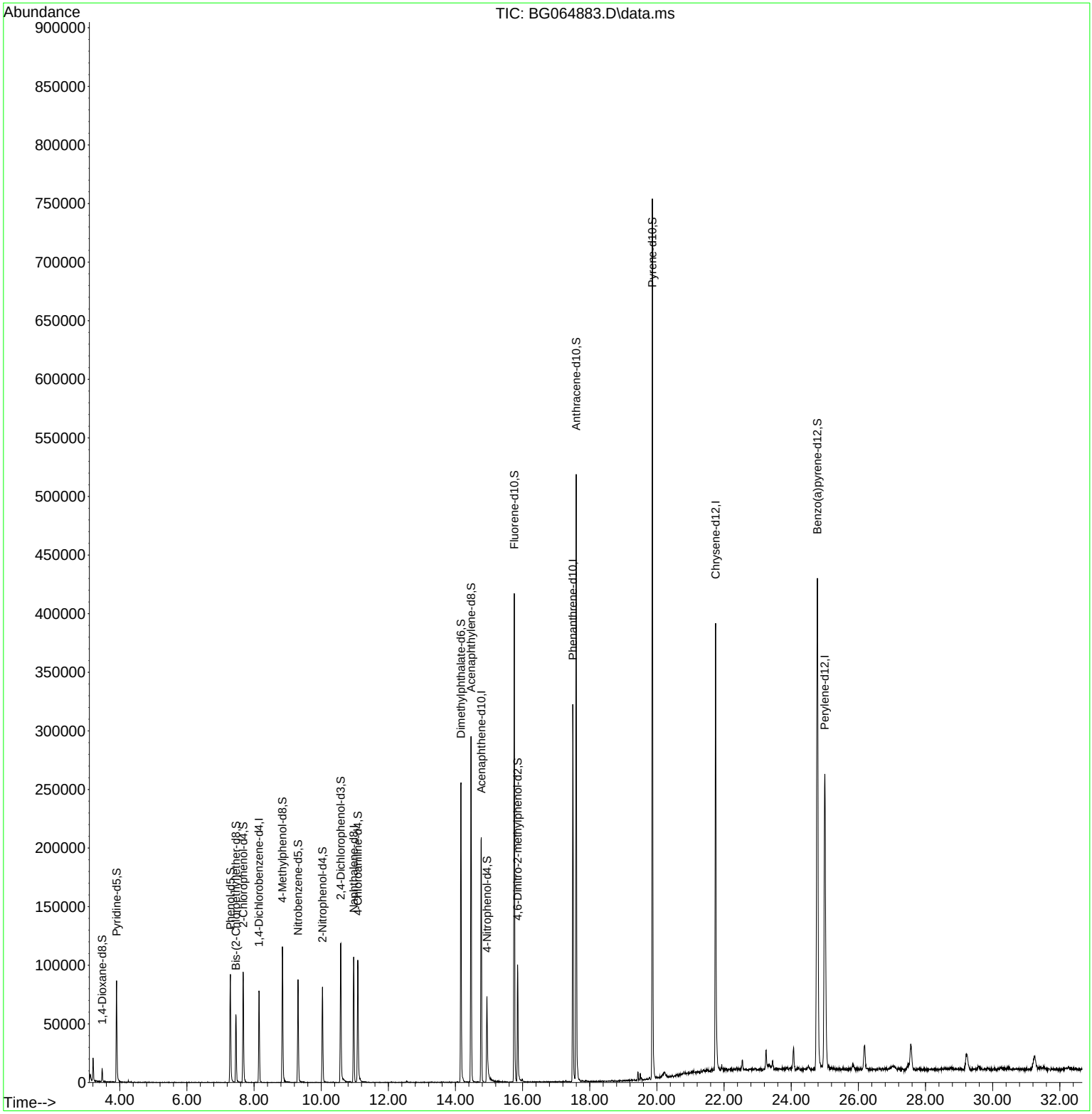
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112525\
Data File : BG064883.D
Acq On : 26 Nov 2025 10:54
Operator : RC/JU
Sample : PB170721BL
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK721

Quant Time: Dec 01 07:50:39 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112425.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 25 10:23:09 2025
Response via : Initial Calibration



Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112525\
Data File : BG064883.D
Acq On : 26 Nov 2025 10:54
Operator : RC/JU
Sample : PB170721BL
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_G
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Quant Time: Dec 01 07:50:39 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112425.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 25 10:23:09 2025
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.147	152	23668	20.000	ng/ul	0.00
20) Naphthalene-d8	10.967	136	91326	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.768	164	86159	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.495	188	223060	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.749	240	299116	20.000	ng/ul	# 0.00
88) Perylene-d12	24.998	264	327094	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.476	96	5427	9.088	ng/uL	0.00
4) Pyridine-d5	3.905	84	40763	24.775	ng/ul	0.00
7) Phenol-d5	7.295	99	53011	27.784	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.459	67	29252	23.972	ng/ul	0.00
11) 2-Chlorophenol-d4	7.677	132	45782	31.652	ng/ul	0.00
15) 4-Methylphenol-d8	8.846	113	45212	27.494	ng/ul	0.00
21) Nitrobenzene-d5	9.310	128	20810	30.211	ng/ul	0.00
24) 2-Nitrophenol-d4	10.039	143	25974	31.747	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.579	165	52716	31.251	ng/ul	0.00
31) 4-Chloroaniline-d4	11.090	131	61048	28.313	ng/ul	0.00
46) Dimethylphthalate-d6	14.163	166	208277	30.731	ng/ul	0.00
49) Acenaphthylene-d8	14.463	160	217238	29.930	ng/ul	0.00
54) 4-Nitrophenol-d4	14.939	143	24797	24.133	ng/ul	0.00
60) Fluorene-d10	15.750	176	187286	30.782	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.850	200	32422	22.703	ng/ul	0.00
73) Anthracene-d10	17.595	188	351131	31.153	ng/ul	0.00
81) Pyrene-d10	19.863	212	482927	31.432	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.780	264	557750	32.320	ng/ul	0.00

Target Compounds	Qvalue

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