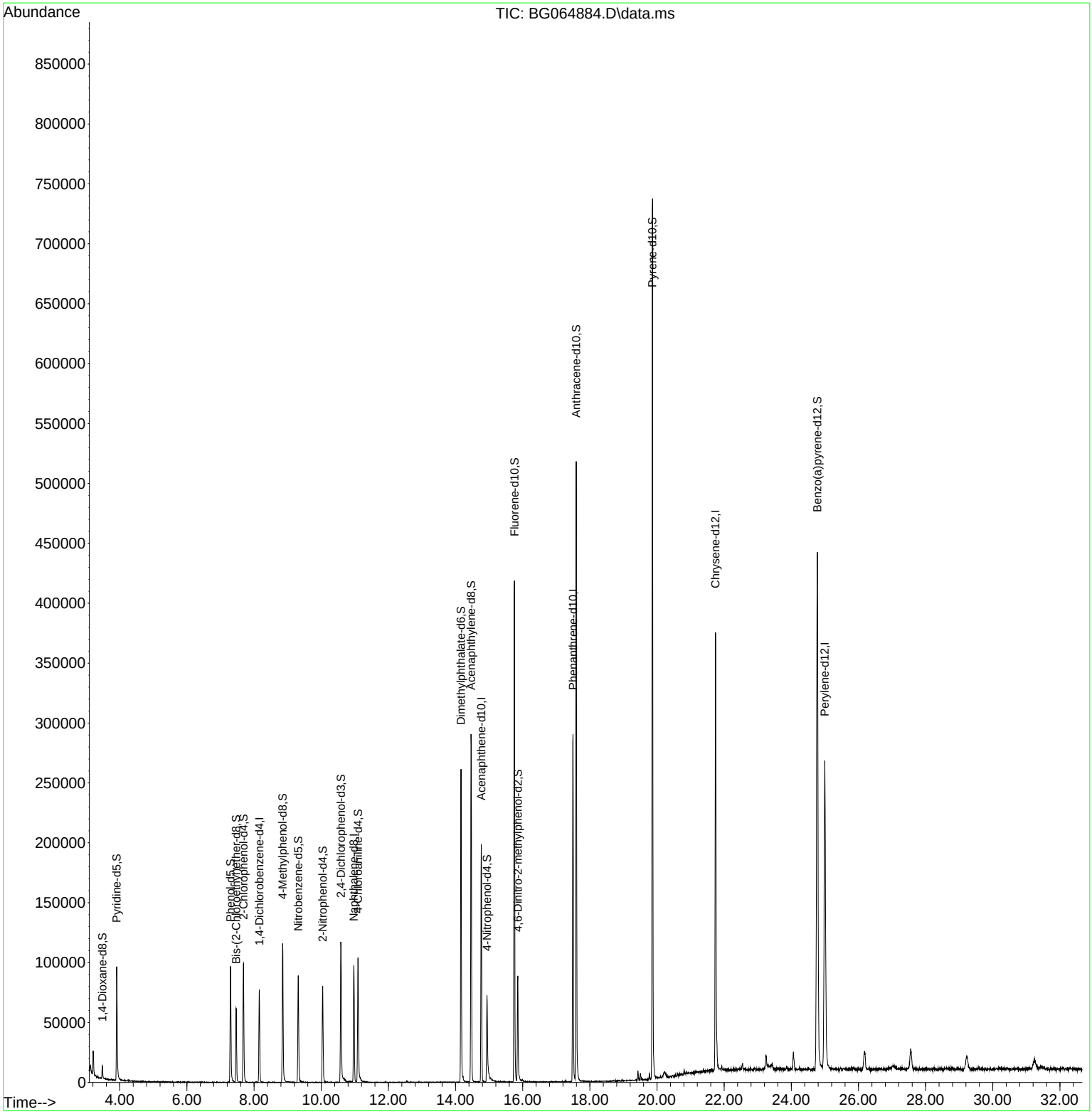
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112525\
Data File : BG064884.D
Acq On : 26 Nov 2025 11:35
Operator : RC/JU
Sample : PB170723BL
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK723

Quant Time: Dec 01 07:51:00 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 : BG064884.D
Acq On : 26 Nov 2025 11:35
Operator : RC/JU
Sample : PB170723BL
Misc :
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Instrument : BNA_G
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Quant Time: Dec 01 07:51:00 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	8.158	152	24283	20.000	ng/ul	0.00
20) Naphthalene-d8	10.972	136	86769	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.768	164	81265	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.500	188	214286	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.742	240	287929	20.000	ng/ul	# 0.00
88) Perylene-d12	24.997	264	319508	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.481	96	5256	8.579	ng/uL	0.00
4) Pyridine-d5	3.910	84	45126	26.732	ng/ul	0.00
7) Phenol-d5	7.300	99	54854	28.022	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.465	67	31635	25.268	ng/ul	0.00
11) 2-Chlorophenol-d4	7.682	132	49364	33.264	ng/ul	0.00
15) 4-Methylphenol-d8	8.851	113	46924	27.813	ng/ul	0.00
21) Nitrobenzene-d5	9.315	128	22109	33.782	ng/ul	0.00
24) 2-Nitrophenol-d4	10.044	143	27638	35.555	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.585	165	52729	32.900	ng/ul	0.00
31) 4-Chloroaniline-d4	11.096	131	62545	30.531	ng/ul	0.00
46) Dimethylphthalate-d6	14.163	166	206594	32.319	ng/ul	0.00
49) Acenaphthylene-d8	14.462	160	223085	32.587	ng/ul	0.00
54) 4-Nitrophenol-d4	14.938	143	22896	23.625	ng/ul	0.00
60) Fluorene-d10	15.755	176	189929	33.096	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.855	200	31604	23.036	ng/ul	0.00
73) Anthracene-d10	17.594	188	362673	33.495	ng/ul	0.00
81) Pyrene-d10	19.862	212	493672	33.379	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.780	264	579229	34.361	ng/ul	0.00
Target Compounds						Qvalue

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