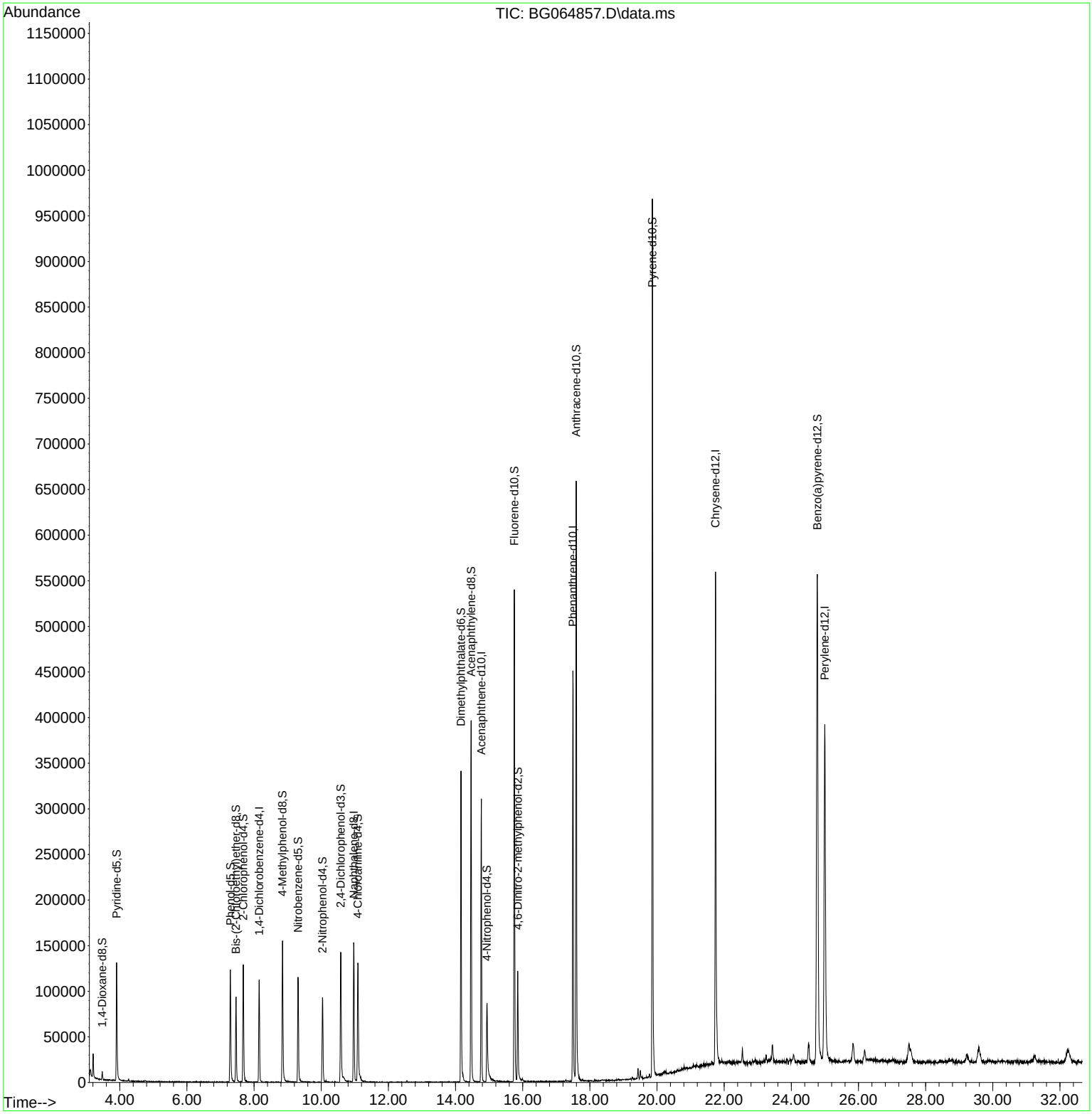
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112525\
Data File : BG064857.D
Acq On : 25 Nov 2025 12:28
Operator : RC/JU
Sample : PB170717BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK717

Quant Time: Nov 25 13:03:30 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 : BG064857.D
Acq On : 25 Nov 2025 12:28
Operator : RC/JU
Sample : PB170717BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SBLK717

Quant Time: Nov 25 13:03:30 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	8.152	152	36472	20.000	ng/ul	0.00
20) Naphthalene-d8	10.966	136	138291	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.768	164	121336	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.494	188	313183	20.000	ng/ul	0.00
79) Chrysene-d12	21.742	240	411186	20.000	ng/ul	0.00
88) Perylene-d12	24.997	264	445430	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.475	96	4345	4.722	ng/uL	0.00
4) Pyridine-d5	3.904	84	65964	26.017	ng/ul	0.00
7) Phenol-d5	7.294	99	78958	26.855	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.459	67	49985	26.582	ng/ul	0.00
11) 2-Chlorophenol-d4	7.676	132	64616	28.990	ng/ul	0.00
15) 4-Methylphenol-d8	8.845	113	65340	25.785	ng/ul	0.00
21) Nitrobenzene-d5	9.309	128	29266	28.058	ng/ul	0.00
24) 2-Nitrophenol-d4	10.038	143	34701	28.009	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.578	165	64672	25.319	ng/ul	0.00
31) 4-Chloroaniline-d4	11.090	131	85360	26.144	ng/ul	0.00
46) Dimethylphthalate-d6	14.162	166	281135	29.455	ng/ul	0.00
49) Acenaphthylene-d8	14.462	160	296434	29.001	ng/ul	0.00
54) 4-Nitrophenol-d4	14.932	143	30219	20.884	ng/ul	0.00
60) Fluorene-d10	15.749	176	250567	29.243	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.855	200	40026	19.962	ng/ul	0.00
73) Anthracene-d10	17.594	188	468300	29.592	ng/ul	0.00
81) Pyrene-d10	19.862	212	638061	30.210	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.774	264	704900	29.995	ng/ul	0.00

Target Compounds	Qvalue

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