

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091225\
 Data File : BG064364.D
 Acq On : 12 Sep 2025 18:08
 Operator : RC/JU
 Sample : PB169653BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK653

Quant Time: Sep 15 10:56:57 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091125.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 12 11:06:20 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.842	152	18536	20.000	ng/u1	0.00
20) Naphthalene-d8	10.633	136	78915	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.470	164	60219	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.208	188	151807	20.000	ng/u1	0.00
79) Chrysene-d12	21.432	240	179872	20.000	ng/u1	# 0.00
88) Perylene-d12	24.435	264	195203	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.318	96	1952	5.532	ng/uL	0.00
4) Pyridine-d5	3.718	84	34780	33.784	ng/u1	0.00
7) Phenol-d5	7.014	99	47018	31.033	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.178	67	24841	31.419	ng/u1	0.00
11) 2-Chlorophenol-d4	7.378	132	38262	31.576	ng/u1	0.00
15) 4-Methylphenol-d8	8.547	113	41368	30.009	ng/u1	0.00
21) Nitrobenzene-d5	8.994	128	19307	31.179	ng/u1	0.00
24) 2-Nitrophenol-d4	9.717	143	24739	32.753	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	41922	29.848	ng/u1	0.00
31) 4-Chloroaniline-d4	10.762	131	55785	29.381	ng/u1	0.00
46) Dimethylphthalate-d6	13.876	166	176914	32.964	ng/u1	0.00
49) Acenaphthylene-d8	14.158	160	161640	31.172	ng/u1	0.00
54) 4-Nitrophenol-d4	14.670	143	21998	26.160	ng/u1	0.00
60) Fluorene-d10	15.463	176	137916	31.673	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.580	200	29198	26.964	ng/u1	0.00
73) Anthracene-d10	17.308	188	247078	32.627	ng/u1	0.00
81) Pyrene-d10	19.593	212	320040	31.588	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.229	264	333459	33.145	ng/u1	-0.01

Target Compounds	Qvalue

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

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