

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091525\
 Data File : BG064378.D
 Acq On : 15 Sep 2025 17:02
 Operator : RC/JU
 Sample : PB169657BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK657

Quant Time: Sep 16 10:50:28 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	20980	20.000	ng/u1	0.00
20) Naphthalene-d8	10.633	136	86240	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.463	164	65342	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.207	188	155195	20.000	ng/u1	0.00
79) Chrysene-d12	21.432	240	187581	20.000	ng/u1	0.00
88) Perylene-d12	24.434	264	206911	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.312	96	2604	6.520	ng/uL	0.00
4) Pyridine-d5	3.717	84	39490	33.891	ng/u1	0.00
7) Phenol-d5	7.013	99	54320	31.676	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.178	67	29335	32.780	ng/u1	0.00
11) 2-Chlorophenol-d4	7.378	132	45404	33.105	ng/u1	0.00
15) 4-Methylphenol-d8	8.547	113	48233	30.913	ng/u1	0.00
21) Nitrobenzene-d5	8.993	128	23046	34.056	ng/u1	0.00
24) 2-Nitrophenol-d4	9.716	143	27812	33.694	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	49790	32.439	ng/u1	0.00
31) 4-Chloroaniline-d4	10.762	131	62964	30.345	ng/u1	0.00
46) Dimethylphthalate-d6	13.876	166	195739	33.612	ng/u1	0.00
49) Acenaphthylene-d8	14.158	160	193049	34.310	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	24715	27.087	ng/u1	0.00
60) Fluorene-d10	15.456	176	161199	34.118	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	31294	28.269	ng/u1	0.00
73) Anthracene-d10	17.307	188	278693	35.999	ng/u1	0.00
81) Pyrene-d10	19.593	212	355328	33.630	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.228	264	388713	36.451	ng/u1	-0.01

Target Compounds	Qvalue

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

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