

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091225\
 Data File : BG064355.D
 Acq On : 12 Sep 2025 12:00
 Operator : RC/JU
 Sample : PB169495BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK495

Quant Time: Sep 12 13:13:34 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	7.847	152	19557	20.000	ng/u1	0.00
20) Naphthalene-d8	10.632	136	83249	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.468	164	61508	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.206	188	157968	20.000	ng/u1	0.00
79) Chrysene-d12	21.437	240	170323	20.000	ng/u1	0.00
88) Perylene-d12	24.433	264	191192	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	2408	6.468	ng/uL	0.00
4) Pyridine-d5	3.716	84	34518	31.779	ng/u1	0.00
7) Phenol-d5	7.012	99	48827	30.544	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.171	67	27398	32.844	ng/u1	0.00
11) 2-Chlorophenol-d4	7.377	132	41140	32.178	ng/u1	0.00
15) 4-Methylphenol-d8	8.546	113	41857	28.779	ng/u1	0.00
21) Nitrobenzene-d5	8.992	128	21179	32.422	ng/u1	0.00
24) 2-Nitrophenol-d4	9.715	143	24815	31.143	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.256	165	45746	30.875	ng/u1	0.00
31) 4-Chloroaniline-d4	10.767	131	59517	29.714	ng/u1	0.00
46) Dimethylphthalate-d6	13.875	166	181429	33.097	ng/u1	0.00
49) Acenaphthylene-d8	14.163	160	170190	32.133	ng/u1	0.00
54) 4-Nitrophenol-d4	14.668	143	23525	27.390	ng/u1	0.00
60) Fluorene-d10	15.461	176	144175	32.417	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.579	200	30799	27.333	ng/u1	0.00
73) Anthracene-d10	17.306	188	252548	32.049	ng/u1	0.00
81) Pyrene-d10	19.598	212	317319	33.076	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.233	264	326468	33.131	ng/u1	0.00

Target Compounds	Qvalue

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

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