

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091225\
 Data File : BG064366.D
 Acq On : 12 Sep 2025 19:30
 Operator : RC/JU
 Sample : PB169655BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK655

Quant Time: Sep 15 11:00:13 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	20150	20.000	ng/u1	0.00
20) Naphthalene-d8	10.632	136	89438	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.468	164	66401	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.206	188	171961	20.000	ng/u1	0.00
79) Chrysene-d12	21.437	240	203716	20.000	ng/u1	0.00
88) Perylene-d12	24.439	264	221462	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	2269	5.915	ng/uL	0.00
4) Pyridine-d5	3.716	84	33697	30.110	ng/u1	0.00
7) Phenol-d5	7.012	99	47506	28.843	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.177	67	25077	29.177	ng/u1	0.00
11) 2-Chlorophenol-d4	7.377	132	39904	30.293	ng/u1	0.00
15) 4-Methylphenol-d8	8.546	113	43004	28.697	ng/u1	0.00
21) Nitrobenzene-d5	8.992	128	20737	29.549	ng/u1	0.00
24) 2-Nitrophenol-d4	9.715	143	24951	29.147	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.256	165	45071	28.315	ng/u1	0.00
31) 4-Chloroaniline-d4	10.767	131	59617	27.705	ng/u1	0.00
46) Dimethylphthalate-d6	13.875	166	191231	32.315	ng/u1	0.00
49) Acenaphthylene-d8	14.157	160	181006	31.657	ng/u1	0.00
54) 4-Nitrophenol-d4	14.662	143	24644	26.578	ng/u1	0.00
60) Fluorene-d10	15.461	176	152744	31.813	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.579	200	30956	25.237	ng/u1	0.00
73) Anthracene-d10	17.306	188	277598	32.361	ng/u1	0.00
81) Pyrene-d10	19.592	212	352524	30.722	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.233	264	376517	32.987	ng/u1	0.00

Target Compounds Qvalue

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

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