

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090925\
 Data File : BP025693.D
 Acq On : 09 Sep 2025 11:35
 Operator : CG/JU
 Sample : PB169495BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK495

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 09/10/2025
 Supervised By :Jagrut Upadhyay 09/10/2025

Quant Time: Sep 09 12:51:37 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090825.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 09 09:26:50 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.725	152	150362	20.000	ng/u1	-0.04
20) Naphthalene-d8	10.501	136	672716	20.000	ng/u1	-0.03
38) Acenaphthene-d10	14.372	164	440537	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.178	188	970314	20.000	ng/u1	0.00
79) Chrysene-d12	21.630	240	1052143	20.000	ng/u1	0.01
88) Perylene-d12	24.995	264	1077339	20.000	ng/u1	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.255	96	24965	6.794	ng/uL	-0.02
4) Pyridine-d5	3.661	84	353337	33.177	ng/u1	-0.03
7) Phenol-d5	6.907	99	380213	31.305	ng/u1	-0.04
9) Bis-(2-Chloroethyl)eth...	7.060	67	242575	32.664	ng/u1	-0.04
11) 2-Chlorophenol-d4	7.266	132	308760	32.636	ng/u1	-0.04
15) 4-Methylphenol-d8	8.431	113	319499	31.771	ng/u1	-0.04
21) Nitrobenzene-d5	8.866	128	150502	29.083	ng/u1	-0.04
24) 2-Nitrophenol-d4	9.584	143	188985	32.129	ng/u1	-0.04
28) 2,4-Dichlorophenol-d3	10.143	165	318162	30.015	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.637	131	443893	29.888	ng/u1	-0.02
46) Dimethylphthalate-d6	13.772	166	1136843	34.135	ng/u1	-0.01
49) Acenaphthylene-d8	14.060	160	1191270	32.575	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.619	143	169048m	30.601	ng/u1	0.02
60) Fluorene-d10	15.378	176	932660	33.628	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.507	200	183455	27.769	ng/u1	-0.01
73) Anthracene-d10	17.283	188	1532725	32.823	ng/u1	0.00
81) Pyrene-d10	19.660	212	1909264	33.156	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.765	264	1892460	34.216	ng/u1	0.04

Target Compounds Qvalue

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

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