

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032025\
 Data File : BP024193.D
 Acq On : 20 Mar 2025 15:33
 Operator : RC/JU
 Sample : PB167169BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK169

Quant Time: Mar 20 16:00:06 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030425.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 04 14:51:41 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.740	152	348093	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.522	136	1460726	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.392	164	965973	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.204	188	1944580	20.000	ng/u1	0.00
79) Chrysene-d12	21.651	240	1873842	20.000	ng/u1	0.00
88) Perylene-d12	25.051	264	1826937	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	63059	7.216	ng/uL	-0.01
4) Pyridine-d5	3.669	84	868397	34.546	ng/u1	-0.01
7) Phenol-d5	6.922	99	1086794	35.189	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.075	67	596214	35.054	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.281	132	846958	35.724	ng/u1	-0.01
15) 4-Methylphenol-d8	8.446	113	871783	35.750	ng/u1	-0.01
21) Nitrobenzene-d5	8.887	128	392505	33.278	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.604	143	419446	34.062	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.146	165	773578	34.938	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.657	131	1139327	33.667	ng/u1	-0.01
46) Dimethylphthalate-d6	13.792	166	2551351	36.449	ng/u1	0.00
49) Acenaphthylene-d8	14.075	160	2758290	33.974	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.604	143	396516	27.669	ng/u1	0.00
60) Fluorene-d10	15.404	176	2101761	35.341	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.528	200	362386	29.516	ng/u1	0.00
73) Anthracene-d10	17.310	188	3329129	35.081	ng/u1	0.01
81) Pyrene-d10	19.675	212	3850752	41.018	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.815	264	3448171	36.406	ng/u1	0.01

Target Compounds Qvalue

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

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