

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012425\
 Data File : BP023710.D
 Acq On : 24 Jan 2025 10:11
 Operator : RC/JU
 Sample : PB166118BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK118

Quant Time: Jan 24 22:31:36 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012225.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 23 00:03:00 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	359838	20.000	ng/u1	0.00
20) Naphthalene-d8	10.575	136	1381656	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.416	164	857276	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.216	188	1814157	20.000	ng/u1	0.00
79) Chrysene-d12	21.651	240	1828866	20.000	ng/u1	0.00
88) Perylene-d12	25.033	264	1979151	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	74997	8.395	ng/uL	0.00
4) Pyridine-d5	3.705	84	1053330	40.857	ng/u1	0.00
7) Phenol-d5	6.958	99	1231859	40.458	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	689148	40.071	ng/u1	0.00
11) 2-Chlorophenol-d4	7.328	132	986457	42.616	ng/u1	0.00
15) 4-Methylphenol-d8	8.481	113	958103	40.016	ng/u1	0.00
21) Nitrobenzene-d5	8.934	128	430028	41.310	ng/u1	0.00
24) 2-Nitrophenol-d4	9.652	143	391539	41.026	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.187	165	856667	42.705	ng/u1	0.00
31) 4-Chloroaniline-d4	10.699	131	1141590	35.584	ng/u1	0.00
46) Dimethylphthalate-d6	13.816	166	2750423	44.611	ng/u1	0.00
49) Acenaphthylene-d8	14.104	160	3044527	42.668	ng/u1	0.00
54) 4-Nitrophenol-d4	14.593	143	432289	34.896	ng/u1	-0.01
60) Fluorene-d10	15.428	176	2367405	44.858	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.528	200	239488	26.807	ng/u1	0.00
73) Anthracene-d10	17.316	188	3823823	43.904	ng/u1	0.00
81) Pyrene-d10	19.680	212	4444034	47.754	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.804	264	4516426	45.592	ng/u1	0.00

Target Compounds Qvalue

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

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