

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012025\
 Data File : BP023664.D
 Acq On : 20 Jan 2025 21:56
 Operator : RC/JU
 Sample : PB166118BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK118

Quant Time: Jan 20 22:45:49 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jan 20 21:50:54 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	715828	20.000	ng/u1	0.00
20) Naphthalene-d8	10.587	136	2869734	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.428	164	1754932	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.216	188	3472990	20.000	ng/u1	-0.01
79) Chrysene-d12	21.669	240	3376018	20.000	ng/u1	0.00
88) Perylene-d12	25.057	264	3718671	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	113638	5.853	ng/uL	0.00
4) Pyridine-d5	3.717	84	1718092	31.744	ng/u1	0.00
7) Phenol-d5	6.969	99	2037490	33.061	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	1108580	30.707	ng/u1	0.00
11) 2-Chlorophenol-d4	7.340	132	1646455	33.829	ng/u1	0.00
15) 4-Methylphenol-d8	8.493	113	1569132	32.826	ng/u1	0.00
21) Nitrobenzene-d5	8.952	128	731572	33.399	ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	659508	45.871	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	1382548	31.615	ng/u1	0.00
31) 4-Chloroaniline-d4	10.710	131	2049044	29.350	ng/u1	0.00
46) Dimethylphthalate-d6	13.834	166	4121195	30.175	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	4833748	29.751	ng/u1	0.00
54) 4-Nitrophenol-d4	14.604	143	644833	28.404	ng/u1	0.00
60) Fluorene-d10	15.434	176	3569935	30.202	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.534	200	326567	28.899	ng/u1	-0.01
73) Anthracene-d10	17.316	188	5557179	29.642	ng/u1	-0.01
81) Pyrene-d10	19.692	212	6206237	33.184	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.827	264	6372611	31.534	ng/u1	0.02

Target Compounds Qvalue

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

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