

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100424\
 Data File : BP022228.D
 Acq On : 04 Oct 2024 14:21
 Operator : RC/JU
 Sample : PB163610BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK610

Quant Time: Oct 04 23:15:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 04 05:25:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.693	152	139965	20.000	ng/u1	0.00
20) Naphthalene-d8	10.451	136	552274	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.298	164	413714	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.098	188	972600	20.000	ng/u1	0.00
79) Chrysene-d12	21.539	240	1009284	20.000	ng/u1	0.00
88) Perylene-d12	24.821	264	1113877	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	15870	4.441	ng/uL	0.00
4) Pyridine-d5	3.628	84	237680	23.270	ng/u1	0.00
7) Phenol-d5	6.875	99	259416	22.658	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.034	67	155228	22.018	ng/u1	0.00
11) 2-Chlorophenol-d4	7.234	132	204283	24.661	ng/u1	0.00
15) 4-Methylphenol-d8	8.393	113	211360	23.297	ng/u1	0.00
21) Nitrobenzene-d5	8.834	128	101490	24.553	ng/u1	0.00
24) 2-Nitrophenol-d4	9.546	143	115900	24.729	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	233624	22.953	ng/u1	0.00
31) 4-Chloroaniline-d4	10.587	131	278290	22.915	ng/u1	0.00
46) Dimethylphthalate-d6	13.710	166	775850	24.260	ng/u1	0.00
49) Acenaphthylene-d8	13.992	160	837897	24.746	ng/u1	0.00
54) 4-Nitrophenol-d4	14.522	143	86995	15.903	ng/u1	0.00
60) Fluorene-d10	15.310	176	675026	23.862	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.434	200	79897	13.232	ng/u1	0.00
73) Anthracene-d10	17.204	188	1093451	24.082	ng/u1	0.00
81) Pyrene-d10	19.580	212	1325927	25.358	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.574	264	1458604	24.897	ng/u1	0.00

Target Compounds Qvalue

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

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