

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
 Data File : BP019213.D
 Acq On : 02 Jan 2024 18:03
 Operator : MA/JU
 Sample : PB158101BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK101

Quant Time: Jan 02 21:43:00 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 03:48:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	128277	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	445082	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.439	164	264531	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	617214	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	723451	20.000	ng/ul	-0.01
88) Perylene-d12	23.662	264	886625	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	25429	6.770	ng/uL	0.00
4) Pyridine-d5	3.658	84	323201	34.284	ng/ul	0.00
7) Phenol-d5	6.975	99	382285	33.232	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	220054	33.066	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132	310588	33.490	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	288636	32.228	ng/ul	0.00
21) Nitrobenzene-d5	8.963	128	138355	35.960	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	153313	36.871	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	297330	35.411	ng/ul	0.00
31) 4-Chloroaniline-d4	10.740	131	364435	35.286	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	772123	33.045	ng/ul	-0.01
49) Acenaphthylene-d8	14.134	160	890257	35.115	ng/ul	0.00
54) 4-Nitrophenol-d4	14.663	143	124790	31.628	ng/ul	0.00
60) Fluorene-d10	15.434	176	690973	34.231	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.563	200	104662	25.866	ng/ul	-0.01
73) Anthracene-d10	17.292	188	1171171	36.335	ng/ul	-0.01
81) Pyrene-d10	19.539	212	1556072	35.293	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.510	264	1817348	35.660	ng/ul	0.00

Target Compounds Qvalue

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

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