

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010923\
 Data File : BP013412.D
 Acq On : 09 Jan 2023 10:21
 Operator : CG/JU
 Sample : PB150090BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK090

Quant Time: Jan 09 10:48:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 06 01:03:32 2023
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.922	152	393322	20.000 ng/u1	0.00
20) Naphthalene-d8	10.734	136	1590817	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.569	164	1015942	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.328	188	2300639	20.000 ng/u1	0.00
79) Chrysene-d12	21.421	240	2196557	20.000 ng/u1	0.01
88) Perylene-d12	23.892	264	2490061	20.000 ng/u1	0.02
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.311	96	61109	6.039 ng/uL	0.00
4) Pyridine-d5	3.740	84	788443	27.936 ng/u1	0.00
7) Phenol-d5	7.081	99	937523	28.725 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	649045	31.522 ng/u1	0.00
11) 2-Chlorophenol-d4	7.452	132	765670	31.458 ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	727322	29.251 ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	346242	30.568 ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	387528	31.254 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.351	165	727359	29.435 ng/u1	0.00
31) 4-Chloroaniline-d4	10.875	131	995502	28.597 ng/u1	0.00
46) Dimethylphthalate-d6	13.981	166	2490510	33.647 ng/u1	0.00
49) Acenaphthylene-d8	14.263	160	2543811	31.048 ng/u1	0.00
54) 4-Nitrophenol-d4	14.769	143	319051	23.549 ng/u1	0.00
60) Fluorene-d10	15.563	176	2121857	32.454 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	360192	24.691 ng/u1	0.01
73) Anthracene-d10	17.433	188	3401287	33.448 ng/u1	0.01
81) Pyrene-d10	19.663	212	4125789	34.522 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.727	264	4079852	33.356 ng/u1	0.02

Target Compounds	Qvalue

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

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