

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
 Data File : BM043717.D
 Acq On : 03 Jan 2024 05:30
 Operator : MA/JU
 Sample : PB158129BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK129

Quant Time: Jan 03 06:41:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 22:31:57 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	176796	20.000	ng/u1	0.00
20) Naphthalene-d8	10.769	136	765234	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.598	164	459051	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.345	188	991377	20.000	ng/u1	0.00
79) Chrysene-d12	21.527	240	881531	20.000	ng/u1	0.00
88) Perylene-d12	23.945	264	1030429	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	28311	5.204	ng/uL	0.00
4) Pyridine-d5	3.793	84	408922	26.119	ng/u1	0.00
7) Phenol-d5	7.134	99	531677	26.436	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.299	67	317524	24.929	ng/u1	0.00
11) 2-Chlorophenol-d4	7.487	132	402559	26.215	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	415848	26.645	ng/u1	0.01
21) Nitrobenzene-d5	9.134	128	202418	27.340	ng/u1	0.00
24) 2-Nitrophenol-d4	9.851	143	222064	29.029	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	380938	28.209	ng/u1	0.01
31) 4-Chloroaniline-d4	10.916	131	582052	29.297	ng/u1	0.00
46) Dimethylphthalate-d6	14.016	166	1111356	26.894	ng/u1	0.00
49) Acenaphthylene-d8	14.292	160	1397563	28.509	ng/u1	0.00
54) 4-Nitrophenol-d4	14.822	143	246533	31.521	ng/u1	0.03
60) Fluorene-d10	15.586	176	1023520	29.952	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	124753	19.027	ng/u1	0.00
73) Anthracene-d10	17.445	188	1649962	29.676	ng/u1	0.00
81) Pyrene-d10	19.733	212	1830662	26.304	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.792	264	1855284	29.177	ng/u1	0.02

Target Compounds	Qvalue

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

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