

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
 Data File : BM036583.D
 Acq On : 19 Sep 2022 16:57
 Operator : CG/JU
 Sample : PB147658BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK658

Quant Time: Sep 19 22:32:57 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 16 02:39:54 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.028	152	338627	20.000 ng/u1	0.00
20) Naphthalene-d8	10.839	136	1512636	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.651	164	1018234	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.386	188	2121855	20.000 ng/u1	0.00
79) Chrysene-d12	21.545	240	2009445	20.000 ng/u1	-0.01
88) Perylene-d12	23.991	264	2017074	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.393	96	43199	4.782 ng/uL	0.00
4) Pyridine-d5	3.816	84	564596	21.432 ng/u1	0.00
7) Phenol-d5	7.175	99	724864	23.953 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	481481	26.199 ng/u1	0.00
11) 2-Chlorophenol-d4	7.551	132	582299	26.727 ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	577981	25.849 ng/u1	0.00
21) Nitrobenzene-d5	9.192	128	284745	27.812 ng/u1	0.00
24) 2-Nitrophenol-d4	9.916	143	264030	28.791 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	526375	24.767 ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	883419	25.001 ng/u1	0.00
46) Dimethylphthalate-d6	14.063	166	1961644	28.564 ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	2188803	26.870 ng/u1	0.00
54) 4-Nitrophenol-d4	14.827	143	301420	22.629 ng/u1	0.00
60) Fluorene-d10	15.639	176	1695406	27.548 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	181731	21.648 ng/u1	0.00
73) Anthracene-d10	17.486	188	2681725	27.642 ng/u1	0.00
81) Pyrene-d10	19.762	212	3021388	30.563 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.833	264	2862023	29.043 ng/u1	0.00

Target Compounds Qvalue

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

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