

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
 Data File : BM036740.D
 Acq On : 26 Sep 2022 11:50
 Operator : CG/JU
 Sample : PB147814BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK814

Quant Time: Sep 26 23:12:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 26 23:10:22 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.004	152	370338	20.000 ng/u1	0.00
20) Naphthalene-d8	10.816	136	1597959	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.628	164	1002566	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.363	188	1995532	20.000 ng/u1	0.00
79) Chrysene-d12	21.533	240	1483477	20.000 ng/u1	0.00
88) Perylene-d12	23.962	264	1308325	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.387	96	53265	5.392 ng/uL	0.00
4) Pyridine-d5	3.811	84	667259	23.160 ng/u1	0.00
7) Phenol-d5	7.157	99	944741	28.546 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	592879	29.499 ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	750015	31.478 ng/u1	0.00
15) 4-Methylphenol-d8	8.710	113	757429	30.974 ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	367272	33.957 ng/u1	0.00
24) 2-Nitrophenol-d4	9.893	143	356987	36.849 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.434	165	683991	30.465 ng/u1	0.00
31) 4-Chloroaniline-d4	10.951	131	1034181	27.705 ng/u1	0.00
46) Dimethylphthalate-d6	14.039	166	2183968	32.299 ng/u1	0.00
49) Acenaphthylene-d8	14.322	160	2529838	31.542 ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	323928	24.699 ng/u1	0.00
60) Fluorene-d10	15.616	176	1948803	32.161 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.727	200	181378	22.973 ng/u1	0.00
73) Anthracene-d10	17.463	188	3006681	32.953 ng/u1	0.00
81) Pyrene-d10	19.745	212	3125800	42.830 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.809	264	2183606	34.162 ng/u1	0.00

Target Compounds	Qvalue

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

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