

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082222\
 Data File : BM036381.D
 Acq On : 23 Aug 2022 08:34
 Operator : CG/JU
 Sample : PB147176BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK176

Quant Time: Aug 24 00:55:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 23 00:55:01 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	276393	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.581	136	1287577	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.427	164	805950	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.174	188	1528936	20.000	ng/u1	0.00
79) Chrysene-d12	21.362	240	1046345	20.000	ng/u1	0.00
88) Perylene-d12	23.686	264	936408	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	42085	5.803	ng/uL	0.00
4) Pyridine-d5	3.622	84	599298	29.965	ng/u1	0.00
7) Phenol-d5	6.957	99	759087	32.175	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.122	67	510967	34.408	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.310	132	607117	32.752	ng/u1	-0.02
15) 4-Methylphenol-d8	8.504	113	600604	34.095	ng/u1	-0.01
21) Nitrobenzene-d5	8.951	128	316549	32.426	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.669	143	326312	32.953	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.210	165	564828	31.923	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.734	131	920681	33.566	ng/u1	0.00
46) Dimethylphthalate-d6	13.851	166	1894401	37.154	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	2181470	34.303	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.657	143	261300	27.046	ng/u1	0.00
60) Fluorene-d10	15.421	176	1670802	36.263	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	198036	25.194	ng/u1	0.00
73) Anthracene-d10	17.274	188	2403110	35.348	ng/u1	0.00
81) Pyrene-d10	19.568	212	2298145	42.340	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.539	264	1667989	36.330	ng/u1	0.00

Target Compounds Qvalue

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

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