

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082222\
 Data File : BM036386.D
 Acq On : 23 Aug 2022 12:04
 Operator : CG/JU
 Sample : PB147178BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK178

Quant Time: Aug 24 00:57:45 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	7.786	152	226926	20.000	ng/u1	0.00
20) Naphthalene-d8	10.586	136	1064051	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.433	164	668468	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.174	188	1301346	20.000	ng/u1	0.00
79) Chrysene-d12	21.362	240	844759	20.000	ng/u1	0.00
88) Perylene-d12	23.691	264	764753	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.204	96	32812	5.511	ng/uL	0.00
4) Pyridine-d5	3.628	84	499188	30.401	ng/u1	0.00
7) Phenol-d5	6.963	99	597893	30.867	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.128	67	416762	34.182	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	472838	31.069	ng/u1	-0.01
15) 4-Methylphenol-d8	8.510	113	491315	33.971	ng/u1	0.00
21) Nitrobenzene-d5	8.957	128	253374	31.407	ng/u1	0.00
24) 2-Nitrophenol-d4	9.680	143	255125	31.177	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	450526	30.812	ng/u1	0.00
31) 4-Chloroaniline-d4	10.739	131	737721	32.545	ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	1521301	35.973	ng/u1	0.00
49) Acenaphthylene-d8	14.127	160	1748126	33.142	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	188866	23.569	ng/u1	0.00
60) Fluorene-d10	15.427	176	1352204	35.384	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.562	200	146067	21.832	ng/u1	0.00
73) Anthracene-d10	17.274	188	1984283	34.292	ng/u1	0.00
81) Pyrene-d10	19.574	212	1875753	42.805	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.538	264	1304264	34.784	ng/u1	0.00

Target Compounds Qvalue

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

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