

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042122\
 Data File : BM034754.D
 Acq On : 21 Apr 2022 19:18
 Operator : CG/JU
 Sample : PB144260BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK260

Quant Time: Apr 22 01:41:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 05:03:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	175187	20.000	ng/u1	0.00
20) Naphthalene-d8	10.751	136	650036	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.575	164	423873	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.316	188	861417	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.492	240	881566	20.000	ng/u1	# 0.00
88) Perylene-d12	23.892	264	921241	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	22962	5.759	ng/uL	0.00
4) Pyridine-d5	3.793	84	266439	26.922	ng/u1	0.00
7) Phenol-d5	7.104	99	352378	26.631	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	193573	26.086	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	331430	29.595	ng/u1	0.00
15) 4-Methylphenol-d8	8.646	113	305996	28.904	ng/u1	-0.01
21) Nitrobenzene-d5	9.098	128	152491	32.114	ng/u1	0.00
24) 2-Nitrophenol-d4	9.828	143	171527	34.791	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	327353	31.521	ng/u1	0.00
31) 4-Chloroaniline-d4	10.881	131	431357	30.718	ng/u1	0.00
46) Dimethylphthalate-d6	13.986	166	1067772	34.603	ng/u1	0.00
49) Acenaphthylene-d8	14.269	160	1184505	29.848	ng/u1	0.00
54) 4-Nitrophenol-d4	14.751	143	147315	28.439	ng/u1	0.00
60) Fluorene-d10	15.569	176	903647	33.843	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.674	200	136723	27.665	ng/u1	0.00
73) Anthracene-d10	17.416	188	1380300	32.906	ng/u1	0.00
81) Pyrene-d10	19.704	212	1634630	34.069	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.733	264	1641623	35.233	ng/u1	0.00

Target Compounds Qvalue

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

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