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# Proposal for Flavor Les Houches Accord format
Block FCINFO # Program information
  1 SUPERISO # flavor calculator
  2 2.6 # version number
Block FMODESEL # Model selection
  1 1 # Extradimension (UED)
Block SMINPUTS # Standard Model inputs
  1 1.27839951e+02 # alpha_em^(-1)
  2 1.16570000e-05 # G_Fermi
  3 1.17200002e-01 # alpha_s(M_Z)
  4 9.11699982e+01 # m_Z(pole)
  5 4.19999981e+00 # m_b(m_b)
  6 1.72399994e+02 # m_top(pole)
  7 1.77699995e+00 # m_tau(pole)
 24 1.27000000e+00 # m_c(m_c)
Block FMASS # Mass spectrum in GeV
#PDG code mass scheme scale particle
  3 1.05000000e-01 1 2.0e+00 # s
  5 4.68000000e+00 3 0 # b
 211 1.39600000e-01 0 0 # pi+
 313 8.91700000e-01 0 0 # K*
 321 4.93700000e-01 0 0 # K+
 421 1.86484000e+00 0 0 # D0
 431 1.96849000e+00 0 0 # D_s+
 521 5.27950000e+00 0 0 # B+
 531 5.36630000e+00 0 0 # B_s
Block FLIFE # Lifetime in sec
#PDG code lifetime particle
 211 2.60330000e-08 # pi+
 321 1.23800000e-08 # K+
 431 5.00000000e-13 # D_s+
 521 1.63800000e-12 # B+
 531 1.42500000e-12 # B_s
Block FCONST # Decay constant in GeV
#PDG code decay constant particle
 431 2.41000000e-01 # D_s+
 521 2.00000000e-01 # B+
 531 2.45000000e-01 # B_s
Block FCONSTRATIO # Ratio of decay constant
#PDG code1 code2 ratio comment
 321 211 1.18900000e+00 # f_K/f_pi
Block FBAG # Bag parameters
#PDG code B-parameter particle
 511 1.26709794e+00 # B_d
 531 1.23000000e+00 # B_s
Block FFORM # Form Factors in GeV
#number value name
  1 4.60000000e-01 # Delta(w) in B->D l nu
  2 1.026e+00 # G(1) in B->D l nu
  3 1.170e+00 # rho^2 in B->D l nu
  4 3.1e-01 # T1(B->K*)
Block FSHAPE # Shape factors
#number value name
  1 5.80000000e-01 # C (b->s gamma)
Block FWCOEF Q= 1.60846e+02 M= 2
#Effective Wilson coefficients in the standard basis
#order number value
  0 2 1.00000000e+00
  0 7 -1.82057567e-01
  0 8 -1.06651571e-01
  1 1 2.33177662e+01
  1 4 5.29677461e-01
  1 7 1.35373179e-01

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 1   8   -6.94496405e-01
 2   1   3.08498153e+02
 2   2   4.91587899e+01
 2   3   -7.01872509e+00
 2   4   1.25624440e+01
 2   5   8.76122785e-01
 2   6   1.64273022e+00
 2   7   7.05439463e-01
 2   8   -4.65529650e+00
Block FWCOEF Q= 2.34384e+00 M= 2
#Effective Wilson coefficients in the standard basis
#order  number  value
 0     1   -8.47809531e-01
 0     2   1.06562816e+00
 0     3   -1.34214747e-02
 0     4   -1.29110603e-01
 0     5   1.36343067e-03
 0     6   2.88022278e-03
 0     7   -3.73787589e-01
 0     8   -1.80398551e-01
 1     1   1.52422776e+01
 1     2   -2.13433897e+00
 1     3   9.52880033e-02
 1     4   -4.81776851e-01
 1     5   -2.10727176e-02
 1     6   -1.22929476e-02
 1     7   2.14544819e+00
 1     8   -5.16870265e-01
 2     7   1.98785400e+01
Block FOBS # Flavor observables
# ParentPDG type value          NDA  ID1 ID2 ID3 ...
 5     1   2.97350499e-04    2    3  22   # BR(b->s gamma)
521    4   8.25882011e-02    2   313  22   # Delta0(B->K* gamma)
531    1   3.46978963e-09     2    13 -13   # BR(B_s->mu+ mu-)
521    1   1.09699841e-04     2   -15  16   # BR(B_u->tau nu)
521    2   9.96640362e-01     2   -15  16   # R(B_u->tau nu)
431    1   4.81251996e-02     2   -15  16   # BR(D_s->tau nu)
431    1   4.96947301e-03     2   -13  14   # BR(D_s->mu nu)
521    1   6.96556180e-03     3   421 -15  16   # BR(B+->D0 tau nu)
521   11   2.97261612e-01     3   421 -15  16   # BR(B+->D0 tau nu)/BR(B+-> D0 e nu)
321   11   6.45414388e-01     2   -13  14   # BR(K->mu nu)/BR(pi->mu nu)
321   12   9.99985822e-01     2   -13  14   # R_l23
Block FOBSERR # Theoretical error for flavor observables at 68% C.L.
# ParentPDG type -ERR      +ERR      NDA  ID1 ID2 ID3 ...
 5     1   0.3e-04  0.3e-04  2    3  22   # BR(b->s gamma)
Block FOBSSM # SM prediction for flavor observables
# ParentPDG type value          NDA  ID1 ID2 ID3 ...
 5     1   2.97350499e-04    2    3  22   # BR(b->s gamma)

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