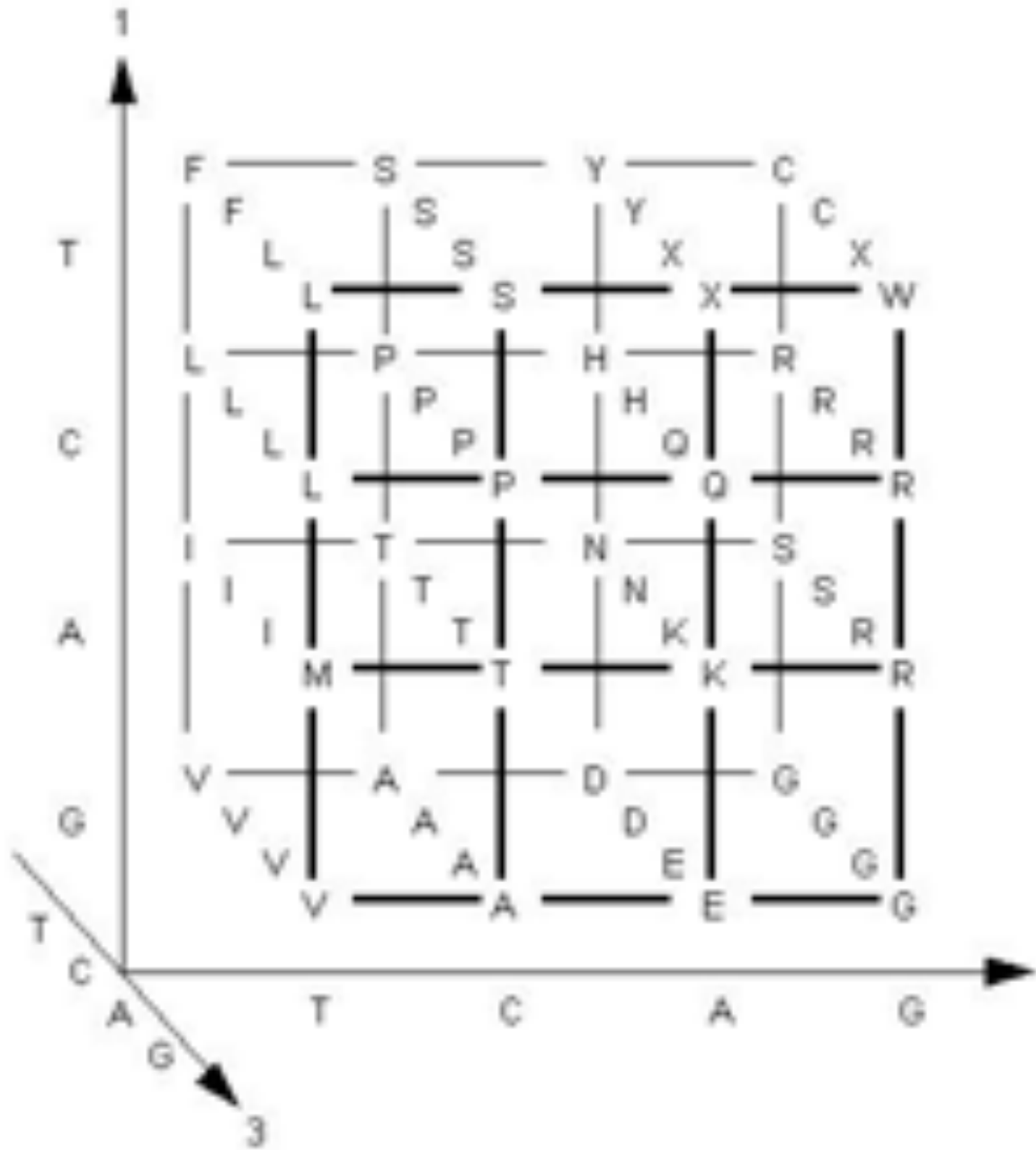


(* Genetic code



*)

(* 64 {a,c,g,t} codons in binary in a,c,g,t order as defined by the following *)

```
t=.; g=.; c=.; a=.;
nucgc = Tuples[{a, c, g, t}, 3];
a = {1, 0, 0, 0};
c = {0, 1, 0, 0};
g = {0, 0, 1, 0};
t = {0, 0, 0, 1};

nucgc = Flatten[nucgc];
nucgc = Partition[nucgc, 12];

pivns = PseudoInverse[nucgc]; (*corrected*)
```

(*

codon triplets entered manually in canonical order by alphabet -
as above in matrix

AAA
AAC
AAG
AAT
ACA
ACC
ACG
ACT
AGA
AGC
AGG
AGT
ATA
ATC
ATG
ATT
CAA
CAC
CAG
CAT
CCA
CCC
CCG
CCT
CGA
CGC
CGG
CGT
CTA
CTC

CTG
CTT
GAA
GAC
GAG
GAT
GCA
GCC
GCG
GCT
GGA
GGC
GGG
GGT
GTA
GTC
GTG
GTT
TAA
TAC
TAG
TAT
TCA
TCC
TCG
TCT
TGA
TGC
TGG
TGT
TTA
TTC
TTG
TTT

manually copy the triplets from above and replace
with amino acid single letter code, in identical order,
and make matrix 'aagc' for amino acid genetic code, including 3 stop codons

*)

```
(* Map an attribute of the amino acids onto the triplet codons *)
```

```
jjt=.
```

```
a=.;
```

```
c=.;
```

```
d=.;
```

```
e=.;
```

```
f=.;
```

```
g=.;
```

```
h=.;
```

```
i=.;
```

```
k=.;
```

```
l=.;
```

```
m=.;
```

```
n=.;
```

```
p=.;
```

```
q=.;
```

```
r=.;
```

```
s=.;
```

```
t=.;
```

```
v=.;
```

```
w=.;
```

```
x=.; y=.;
```

```
jjt = {{k}, {n}, {k}, {n}, {t}, {t}, {t}, {t}, {r}, {s}, {r}, {s}, {i}, {i}, {m}, {i},  
       {q}, {h}, {q}, {h}, {p}, {p}, {p}, {p}, {r}, {r}, {r}, {r}, {l}, {l}, {l}, {l},  
       {e}, {d}, {e}, {d}, {a}, {a}, {a}, {a}, {g}, {g}, {g}, {g}, {v}, {v}, {v}, {v},  
       {x}, {y}, {x}, {y}, {s}, {s}, {s}, {s}, {x}, {c}, {w}, {c}, {l}, {f}, {l}, {f}};
```

```
nuchyd=.;
```

```
nuchyd = pivns.jjt;
```

```

(* 20 K&D hydropathy values, padded with 0.0 for stop *)
(* checked against http://
  web.expasy.org/protscale/pscale/Hphob.Doolittle.html *)

kd = {1.8, 2.5, -3.5001, -3.5002, 2.8, -0.4, -3.2, 4.5, -3.9, 3.8, 1.9,
      -3.5003, -1.6, -3.5004, -4.5, -0.8, -0.7, 4.2, -0.9, 0.0, -1.3};

a = kd[[1]];
c = kd[[2]];
d = kd[[3]];
e = kd[[4]];
f = kd[[5]];
g = kd[[6]];
h = kd[[7]];
i = kd[[8]];
k = kd[[9]];
l = kd[[10]];
m = kd[[11]];
n = kd[[12]];
p = kd[[13]];
q = kd[[14]];
r = kd[[15]];
s = kd[[16]];
t = kd[[17]];
v = kd[[18]];
w = kd[[19]];
x = kd[[20]];
y = kd[[21]];

nuchyd (* numeric with above hydropathy values *)

{{-0.638625}, {-1.25125}, {0.68625}, {0.880375}, {-2.63938}, {-0.163375},
  {-1.46963}, {3.94913}, {-0.169937}, {0.117688}, {-0.388687}, {0.117688}}

(* Reconstruct hydropathy values from nucleotide information
  using dot product of a weighted vector for triplets *)
(* use single letter code for new vectors, precede by the letter 'a',
  because variables are global *)

(* Watch out! Triplets are in A,C,G,T for use here *)

aa = {0, 0, 1, 0, 0, 1, 0, 0, .25, .25, .25, .25};
aanuc = Total[aa.nuchyd, 2]

ac = {0, 0, 0, 1, 0, 0, 1, 0, 0, 0.5, 0, 0.5};

```

```

acnuc = Total[ac.nuchyd, 2]

ad = {0, 0, 1, 0, 1, 0, 0, 0, 0, 0.5, 0.0, 0.5};
adnuc = Total[ad.nuchyd, 2]

ae = {0, 0, 1, 0, 1, 0, 0, 0, 0.5, 0, 0.5, 0};
aenuc = Total[ae.nuchyd, 2]

af = {0, 0, 0, 1, 0, 0, 0, 1, 0, 1, 0, 1};
afnuc = Total[af.nuchyd, 2]

ag = {0, 0, 1, 0, 0, 0, 1, 0, .25, .25, .25, .25};
agnuc = Total[ag.nuchyd, 2]

ah = {0, 1, 0, 0, 1, 0, 0, 0, 0, 0.5, 0, 0.5};
ahnuc = Total[ah.nuchyd, 2]

ai = {1, 0, 0, 0, 0, 0, 0, 1, 1/3, 1/3, 0, 1/3};
ainuc = Total[ai.nuchyd, 2]

ak = {1, 0, 0, 0, 1, 0, 0, 0, 0.5, 0, 0.5, 0};
aknuc = Total[ak.nuchyd, 2]

al = {0, 2/3, 0, 1/3, 0, 0, 0, 1, 2/6, 1/6, 2/6, 1/6};
alnuc = Total[al.nuchyd, 2]

am = {1, 0, 0, 0, 0, 0, 0, 1, 0, 0, 1, 0};
amnuc = Total[am.nuchyd, 2]

an = {1, 0, 0, 0, 1, 0, 0, 0, 0, 0.5, 0, 0.5};
annuc = Total[an.nuchyd, 2]

ap = {0, 1, 0, 0, 0, 1, 0, 0, .25, .25, .25, .25};
apnuc = Total[ap.nuchyd, 2]

aq = {0, 1, 0, 0, 1, 0, 0, 0, 0.5, 0, 0.5, 0};
aqnuc = Total[aq.nuchyd, 2]

ar = {1/3, 2/3, 0, 0, 0, 0, 1, 0, 1/3, 1/6, 1/3, 1/6};
arnuc = Total[ar.nuchyd, 2]

as = {1/3, 0, 0, 2/3, 0, 2/3, 1/3, 0, 1/6, 1/3, 1/6, 1/3};
asnuc = Total[as.nuchyd, 2]

at = {1, 0, 0, 0, 0, 1, 0, 0, .25, .25, .25, .25};
atnuc = Total[at.nuchyd, 2]

av = {0, 0, 1, 0, 0, 0, 0, 1, .25, .25, .25, .25};
avnuc = Total[av.nuchyd, 2]

```

```
aw = {0, 0, 0, 1, 0, 0, 1, 0, 0, 0, 1, 0};  
awnuc = Total[aw.nuchyd, 2]
```

```
ax = {0, 0, 0, 1, 2/3, 0, 1/3, 0, 2/3, 0, 1/3, 0};  
axnuc = Total[ax.nuchyd, 2]
```

```
ay = {0, 0, 0, 1, 1, 0, 0, 0, 0, 1/2, 0, 1/2};  
aynuc = Total[ay.nuchyd, 2]
```

0.442063

-0.471563

-1.83544

-2.23244

5.06488

-0.864187

-3.77294

3.33231

-3.55731

3.26144

2.92181

-3.16031

-1.49544

-4.16994

-2.66365

-0.239396

-0.882813

4.55456

-0.977938

-1.61194

-1.64131

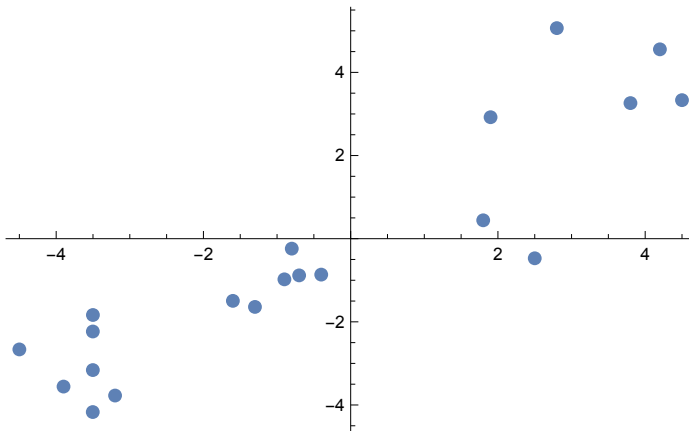
```
(* Fill empirical and SVP values for hydropathy into a 20 x 2 matrix *)
```

```
paired = {{a, aanuc}, {c, acnuc}, {d, adnuc}, {e, aenuc},
  {f, afnuc}, {g, agnuc}, {h, ahnuc}, {i, ainuc}, {k, aknuc},
  {l, alnuc}, {m, amnuc}, {n, annuc}, {p, apnuc}, {q, aqnuc}, {r, arnuc},
  {s, asnuc}, {t, atnuc}, {v, avnuc}, {w, awnuc}, {y, aynuc}};
```

```
ListPlot[paired, PlotStyle -> PointSize[0.02]]
```

```
FindFit[paired, intercept + slope * xaxis, {intercept, slope}, xaxis]
```

```
Correlation[Take[paired, All, {1}], Take[paired, All, {2}]]
```



```
{intercept -> 0.00335099, slope -> 0.862013}
```

```
{{0.914202}}
```


(*
Reaction Centers of Photosynthetic Bacteria (1990) Michel-Beyerle M.ed.pp.209-218,
Springer-Verlag Berlin. Genetic Coding Algorithms for
Engineering Membrane Proteins. Yang M.M., Coleman, W.J., & Youvan,
D.C. Massachusetts Institute of Technology, Department of Chemistry,
Cambridge 02139. Abstract available at
[http://
www.kairos-scientific.com/searchable/abstracts/Yang1990.htm](http://www.kairos-scientific.com/searchable/abstracts/Yang1990.htm) and quoted here:

A solution to the problem of relating the
physicochemical properties of the amino acids to their codon
sequences has been achieved by treating the genetic code as a
system of linear equations and applying the numerical method,
Singular Value Decomposition (SVD). For example,
hydropathy
and molar volume,
which are important determinants of protein
structure and function,
can be quantitatively related to the nucleotide sequence. The
20 hydropathy values of the amino acid residues were
remapped to 12 nucleotide-determined values which,
in turn, were used to predict structural aspects the
photosynthetic reaction center protein,
without DNA→protein translation.

*)

```
(* hotmap *)
```

```
hotmap = Table[{
  Piecewise[{
    {1, 1 <= i < 42},
    {1, 42 <= i < 84},
    {Floor[5.8 * (i - 83)], 84 <= i < 128},
    {256, 128 <= i < 170},
    {256, 170 <= i < 212},
    {256, 212 <= i <= 256}
  ]},
  Piecewise[{
    {1, 1 <= i < 42},
    {(i - 41) * 6, 42 <= i < 84},
    {256, 84 <= i < 128},
    {256 - (i - 128) * 6, 128 <= i < 170},
    {1, 170 <= i < 212},
    {Floor[(i - 211) * 5.7], 212 <= i <= 256}
  ]},
  Piecewise[{
    {Floor[i * 5.8], 1 <= i < 42},
    {256 - ((i - 42) * 6), 42 <= i < 84},
    {1, 84 <= i < 128},
    {1, 128 <= i < 170},
    {(i - 169) * 6, 170 <= i < 212},
    {256, 212 <= i <= 256}
  ]}
],
{i, 1, 256, 1}];
```

```
hotmap[[1, All]] = 1; (* This changes 1,1,
5 to 1,1,1 at i=1 for the blue channel *)
```

```
(* Pseudocolor 61 x 12 pivns matrix *)
```

```
(* get scaling factors *)
```

```
graymat = pivns;
graymin = Min[graymat]
graymax = Max[graymat]
delta = graymax - graymin
```

$$-\frac{1}{96}$$

$$\frac{5}{96}$$

$$\frac{1}{16}$$

```
(* rescale graymat according to min max to get grayscales of 1 to 256 *)

For[x = 1, x ≤ 12, x++,
  For[y = 1, y ≤ 61, y++,
    graymat[[x, y]] = Floor[(graymat[[x, y]] - (graymin)) * (255 / delta) + 1] ]];

colormat = Table[{x, y}, {x, 12}, {y, 61}];
colormat[[All, All]] = {0, 0, 0};
For[x = 1, x ≤ 12, x++,
  For[y = 1, y ≤ 61, y++,
    colormat[[x, y]] = hotmap[[graymat[[x, y]]]] / 256
  ]];

gmat = Graphics[RasterArray[Apply[RGBColor, colormat, {2}]],
  ImageSize → {10 * 61, 10 * 12}, AspectRatio → Automatic];
Show[
  gmat]
```

Part::partw : Part 17387547842683 of

```
{{1, 1, 1}, {1, 1, 11}, {1, 1, 17}, {1, 1, 23}, {1, 1, 29}, {1, 1, 34}, {1, 1, 40}, {1, 1, 46}, {1, 1, 52}, {1, 1, 58}, {1, 1, 63}, {1, 1, 69}, {1, 1, 75},
{1, 1, 81}, {1, 1, 87}, {1, 1, 92}, {1, 1, 98}, {1, 1, 104}, {1, 1, 110}, {1, 1, 116}, {1, 1, 121}, <<9>>, {1, 1, 179}, {1, 1, 185}, {1, 1, 191},
{1, 1, 197}, {1, 1, 203}, {1, 1, 208}, {1, 1, 214}, {1, 1, 220}, {1, 1, 226}, {1, 1, 232}, {1, 1, 237}, {1, 6, 256}, {1, 12, 250}, {1, 18, 244},
{1, 24, 238}, {1, 30, 232}, {1, 36, 226}, {1, 42, 220}, {1, 48, 214}, {1, 54, 208}, <<206>>} does not exist. >>
```

Part::partw : Part 17387547842683 of

```
{{1, 1, 1}, {1, 1, 11}, {1, 1, 17}, {1, 1, 23}, {1, 1, 29}, {1, 1, 34}, {1, 1, 40}, {1, 1, 46}, {1, 1, 52}, {1, 1, 58}, {1, 1, 63}, {1, 1, 69}, {1, 1, 75},
{1, 1, 81}, {1, 1, 87}, {1, 1, 92}, {1, 1, 98}, {1, 1, 104}, {1, 1, 110}, {1, 1, 116}, {1, 1, 121}, <<9>>, {1, 1, 179}, {1, 1, 185}, {1, 1, 191},
{1, 1, 197}, {1, 1, 203}, {1, 1, 208}, {1, 1, 214}, {1, 1, 220}, {1, 1, 226}, {1, 1, 232}, {1, 1, 237}, {1, 6, 256}, {1, 12, 250}, {1, 18, 244},
{1, 24, 238}, {1, 30, 232}, {1, 36, 226}, {1, 42, 220}, {1, 48, 214}, {1, 54, 208}, <<206>>} does not exist. >>
```

Part::partw : Part 17387547842683 of

```
{{1, 1, 1}, {1, 1, 11}, {1, 1, 17}, {1, 1, 23}, {1, 1, 29}, {1, 1, 34}, {1, 1, 40}, {1, 1, 46}, {1, 1, 52}, {1, 1, 58}, {1, 1, 63}, {1, 1, 69}, {1, 1, 75},
{1, 1, 81}, {1, 1, 87}, {1, 1, 92}, {1, 1, 98}, {1, 1, 104}, {1, 1, 110}, {1, 1, 116}, {1, 1, 121}, <<9>>, {1, 1, 179}, {1, 1, 185}, {1, 1, 191},
{1, 1, 197}, {1, 1, 203}, {1, 1, 208}, {1, 1, 214}, {1, 1, 220}, {1, 1, 226}, {1, 1, 232}, {1, 1, 237}, {1, 6, 256}, {1, 12, 250}, {1, 18, 244},
{1, 24, 238}, {1, 30, 232}, {1, 36, 226}, {1, 42, 220}, {1, 48, 214}, {1, 54, 208}, <<206>>} does not exist. >>
```

General::stop : Further output of Part::partw will be suppressed during this calculation. >>

RasterArray::obs : RasterArray is obsolete. Translating to Raster. >>

