beta.dat

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#AA CA      CB       CO     HN    HA       NH
A 50.990 21.730 175.883 8.625 4.923 124.525
C 57.181 30.349 173.744 8.748 5.010 121.913
D 53.618 42.519 175.671 8.536 4.943 122.889
E 55.083 32.343 175.382 8.570 4.846 122.255
F 56.500 41.661 174.563 8.799 5.120 121.385
G 45.048 00.000 172.632 8.350 0.000 109.579
H 54.923 32.261 174.213 8.671 5.048 121.072
I 59.972 40.011 175.039 8.738 4.667 122.884
K 55.153 34.789 175.418 8.554 4.807 122.577
L 53.858 44.247 175.667 8.754 4.872 124.577
M 54.297 35.328 174.926 8.734 5.021 121.709
N 52.378 40.334 174.741 8.630 5.066 121.697
P 62.600 32.160 176.525 0.000 4.595 134.923
Q 54.622 31.422 174.932 8.585 4.810 121.734
R 54.857 32.783 175.018 8.611 4.836 122.233
S 57.210 65.195 173.771 8.568 5.004 117.115
T 61.080 70.679 173.834 8.590 4.929 117.358
V 60.812 33.991 174.910 8.683 4.639 122.302
W 56.268 31.631 175.298 8.740 5.190 122.862
Y 56.689 41.085 174.678 8.772 5.095 121.461
X 54.784 43.342 173.793 8.778 5.207 120.738
#pH 3
D 52.708 41.352 174.935 8.454 5.167 121.134
E 54.911 30.763 175.535 8.615 4.843 122.671
H 54.975 30.275 174.144 8.616 5.102 120.021
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#AA CA      CB       CO     HN    HA       NH
A 52.080 19.167 176.648 8.190 4.295 125.337
C 58.096 29.718 174.634 8.227 4.425 120.362
D 54.127 40.993 176.026 8.209 4.611 121.112
E 56.598 30.334 176.174 8.307 4.262 122.645
F 58.052 39.744 175.391 8.062 4.561 121.150
G 45.165 00.000 175.459 8.165 0.000 113.039
H 55.691 29.529 174.414 8.365 4.624 120.590
I 61.060 38.705 176.011 7.888 4.128 121.160
K 56.219 32.976 176.206 8.142 4.283 123.316
L 54.768 42.270 176.488 8.026 4.339 123.164
M 55.440 32.701 175.359 8.341 4.417 121.452
N 53.040 38.977 174.718 8.288 4.696 120.493
P 62.789 31.824 176.496 0.000 4.425 134.680
Q 55.827 29.614 175.501 8.181 4.284 121.652
R 56.184 30.878 176.142 8.168 4.291 122.938
S 58.380 63.953 174.610 8.313 4.476 117.260
T 61.860 69.810 174.776 8.082 4.324 114.527
V 62.269 32.773 175.848 7.880 4.086 121.171
W 56.740 29.297 175.318 7.955 4.611 121.997
Y 58.024 39.045 175.222 7.984 4.521 121.264
X 54.896 39.833 176.062 8.236 4.727 119.470
#pH 3
D 53.305 38.348 175.033 8.539 4.564 119.648
E 55.841 29.704 175.933 8.197 4.404 120.382
H 55.691 29.529 174.414 8.365 4.624 120.590
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#AA CA      CB       CO     HN    HA       NH
A 54.774 18.223 178.855 8.076 4.025 121.926
C 62.739 27.526 175.925 8.125 4.087 118.292
D 56.941 40.466 177.803 8.207 4.410 119.871
E 59.070 29.376 178.123 8.223 4.017 119.578
F 60.677 39.010 176.592 8.212 4.174 119.930
G 46.817 00.000 175.229 8.249 0.000 107.634
H 58.947 29.787 176.448 8.066 4.304 118.462
I 64.481 37.616 177.119 8.046 3.664 120.080
K 59.010 32.181 177.939 7.977 3.981 119.660
L 57.456 41.629 177.987 8.082 4.006 120.144
M 57.981 32.166 177.440 8.075 4.116 118.786
N 55.478 38.412 176.558 8.183 4.485 117.826
P 65.521 31.476 178.050 0.000 4.202 131.771
Q 58.473 28.444 177.543 8.061 3.993 118.776
R 58.975 29.948 177.775 8.077 3.981 119.487
S 61.114 62.969 175.775 8.151 4.202 115.180
T 65.714 68.651 175.659 8.051 3.989 115.138
V 66.033 31.510 177.149 8.008 3.599 119.520
W 60.181 29.143 177.362 8.228 4.303 120.857
Y 60.947 38.295 176.783 8.151 4.150 119.807
X 58.325 38.764 175.550 8.210 4.348 117.679
#pH 3
D 54.797 39.186 176.723 8.226 4.473 118.391
E 56.968 28.532 176.616 8.136 4.087 120.292
H 55.774 28.672 176.117 8.557 4.403 119.692
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#AA CA      CB       CO     HN     HA       NH
A 52.156 19.321 177.195 8.343 4.341 124.696
C 57.528 30.478 175.189 8.779 4.519 123.546
D 54.263 41.587 175.995 8.500 4.641 122.238
E 56.013 30.492 176.094 8.433 4.374 122.002
F 56.675 40.233 174.779 8.487 4.663 121.914
G 44.548 00.000 173.308 8.478 0.000 110.191
H 55.177 30.454 174.292 8.403 4.701 120.668
I 59.954 38.798 175.794 8.251 4.264 123.115
K 55.804 33.237 176.148 8.341 4.326 122.684
L 54.166 42.594 176.323 8.388 4.435 123.756
M 54.884 33.483 175.085 8.401 4.561 121.512
N 52.899 38.980 174.927 8.542 4.754 120.989
P 62.663 31.925 176.478 0.000 4.414 134.088
Q 55.051 29.569 175.293 8.447 4.403 121.707
R 55.806 30.843 176.052 8.357 4.336 122.030
S 58.329 64.227 174.255 8.539 4.448 117.768
T 61.157 69.996 174.195 8.324 4.448 116.831
V 61.691 32.802 175.557 8.264 4.152 121.839
W 58.358 30.339 176.309 8.491 4.516 123.624
Y 57.982 39.594 175.990 8.637 4.563 121.626
X 55.067 41.106 174.430 8.632 4.632 121.673
#pH 3
D 54.654 38.983 176.289 8.558 4.733 119.932
E 56.708 30.042 176.151 7.995 4.568 121.845
H 55.177 30.454 174.292 8.403 4.701 120.668
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