



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 02:44 pm BST

PDB ID : 1VDM
Title : Crystal structure of purine phosphoribosyltransferase from *Pyrococcus horikoshii* Ot3
Authors : Sugahara, M.; Kunishima, N.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2004-03-23
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

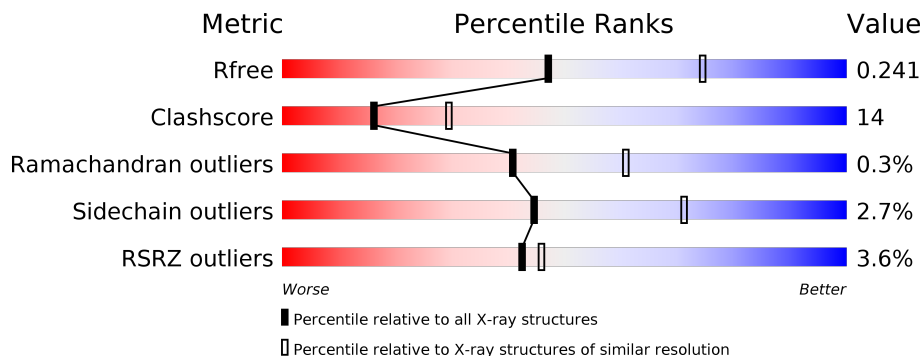
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	153	<div style="display: flex; align-items: center;"> <div style="width: 7%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 62%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 34%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: red;"></div> </div> <p style="text-align: center;">62% 34%</p>
1	B	153	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 64%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 33%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: red;"></div> </div> <p style="text-align: center;">64% 33%</p>
1	C	153	<div style="display: flex; align-items: center;"> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 69%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 24%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 7%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">69% 24% 7%</p>
1	D	153	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 68%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 27%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: red;"></div> </div> <p style="text-align: center;">68% 27%</p>
1	E	153	<div style="display: flex; align-items: center;"> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 66%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 30%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: red;"></div> </div> <p style="text-align: center;">66% 30%</p>
1	F	153	<div style="display: flex; align-items: center;"> <div style="width: 82%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 15%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: red;"></div> </div> <p style="text-align: center;">82% 15%</p>

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Mol	Chain	Length	Quality of chain
1	G	153	<p>2% 64% 33% ..</p>
1	H	153	<p>3% 64% 31% ..</p>
1	I	153	<p>% 71% 24% • 5%</p>
1	J	153	<p>% 75% 20% ..</p>
1	K	153	<p>5% 61% 33% • 5%</p>
1	L	153	<p>5% 67% 30% •</p>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 15181 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called purine phosphoribosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	151	Total 1227	C 810	N 199	O 215	S 3	0	0	0
1	B	149	Total 1207	C 799	N 194	O 211	S 3	0	0	0
1	C	143	Total 1159	C 770	N 187	O 199	S 3	0	0	0
1	D	147	Total 1193	C 792	N 192	O 206	S 3	0	0	0
1	E	149	Total 1211	C 803	N 194	O 211	S 3	0	0	0
1	F	150	Total 1213	C 804	N 195	O 211	S 3	0	0	0
1	G	152	Total 1233	C 815	N 200	O 215	S 3	0	0	0
1	H	149	Total 1214	C 803	N 197	O 211	S 3	0	0	0
1	I	146	Total 1184	C 786	N 190	O 205	S 3	0	0	0
1	J	147	Total 1195	C 792	N 194	O 206	S 3	0	0	0
1	K	146	Total 1189	C 790	N 191	O 205	S 3	0	0	0
1	L	149	Total 1207	C 799	N 194	O 211	S 3	0	0	0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	46	Total 46	O 46	0	0
2	B	69	Total 69	O 69	0	0

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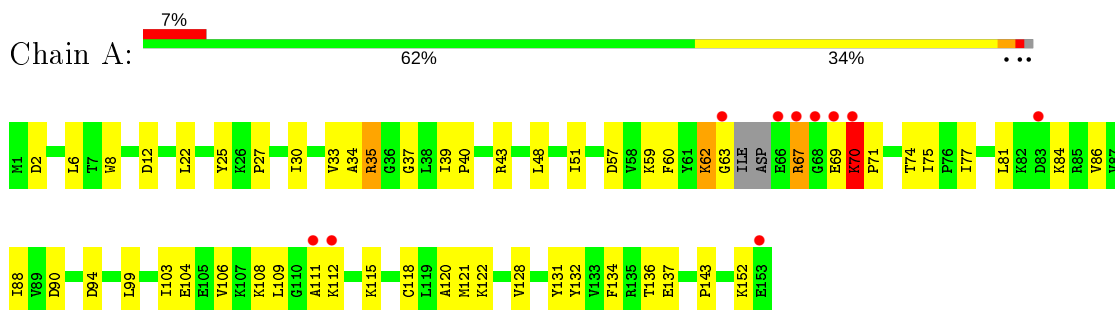
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	52	Total 52	O 52	0	0
2	D	57	Total 57	O 57	0	0
2	E	92	Total 92	O 92	0	0
2	F	98	Total 98	O 98	0	0
2	G	52	Total 52	O 52	0	0
2	H	51	Total 51	O 51	0	0
2	I	52	Total 52	O 52	0	0
2	J	59	Total 59	O 59	0	0
2	K	52	Total 52	O 52	0	0
2	L	69	Total 69	O 69	0	0

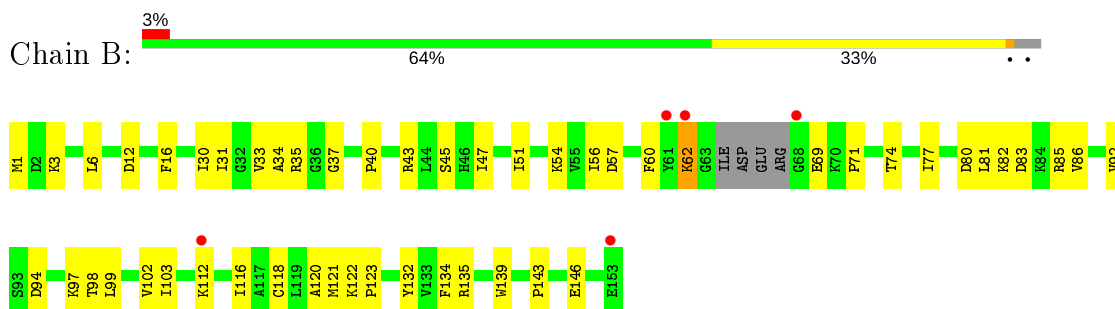
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

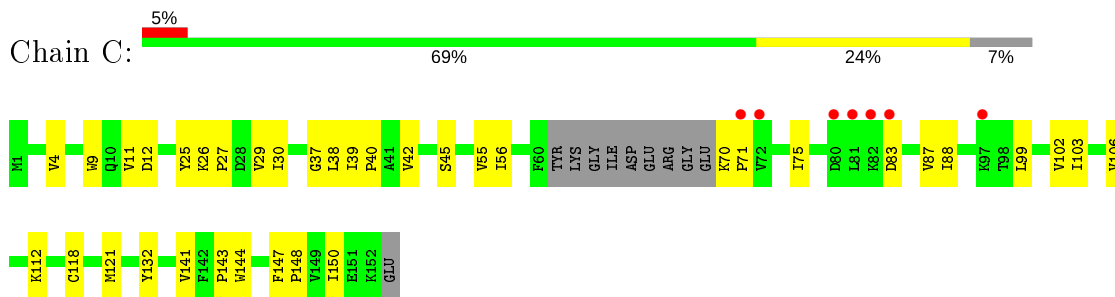
- Molecule 1: purine phosphoribosyltransferase



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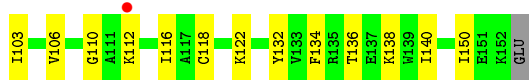
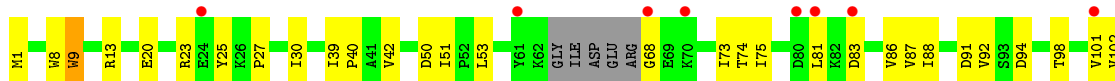


- Molecule 1: purine phosphoribosyltransferase

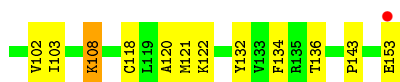
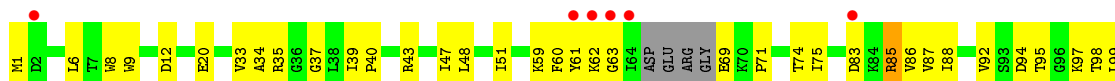


- Molecule 1: purine phosphoribosyltransferase

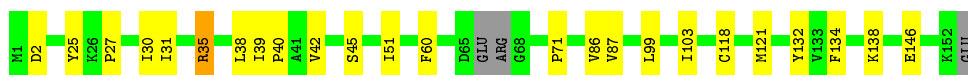
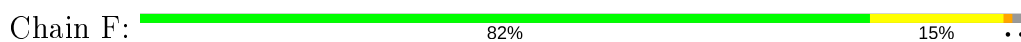




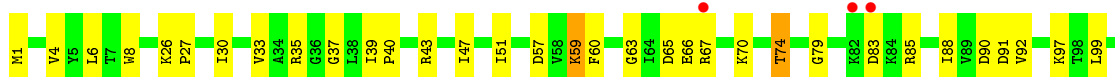
- Molecule 1: purine phosphoribosyltransferase



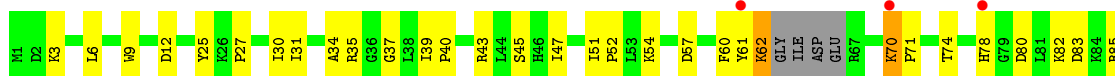
- Molecule 1: purine phosphoribosyltransferase



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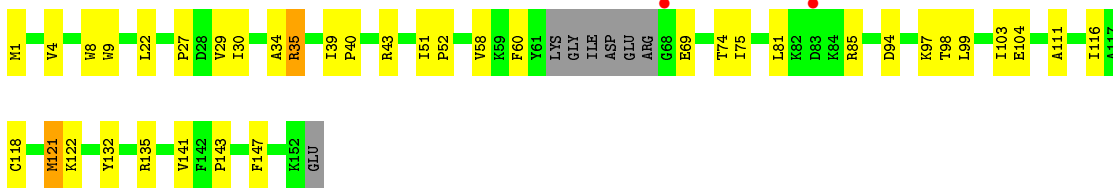


- Molecule 1: purine phosphoribosyltransferase

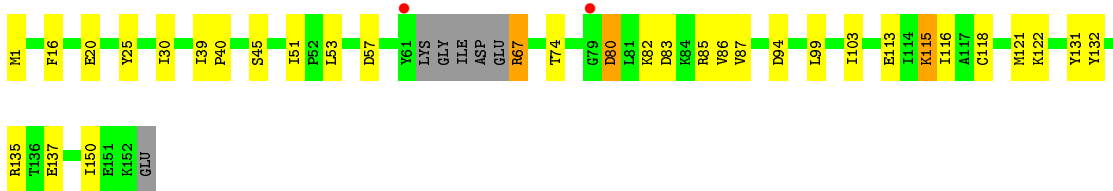
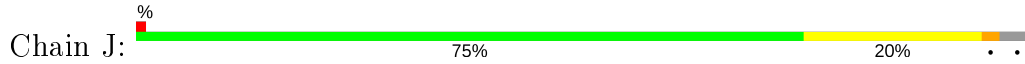


- Molecule 1: purine phosphoribosyltransferase

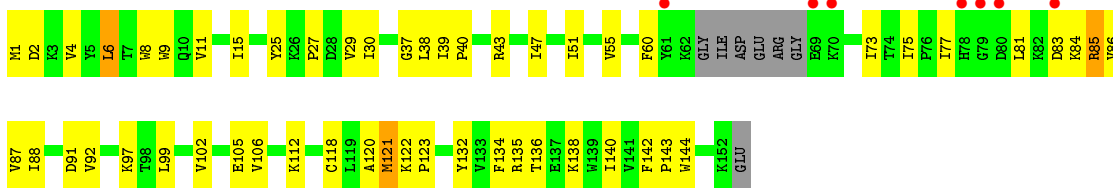




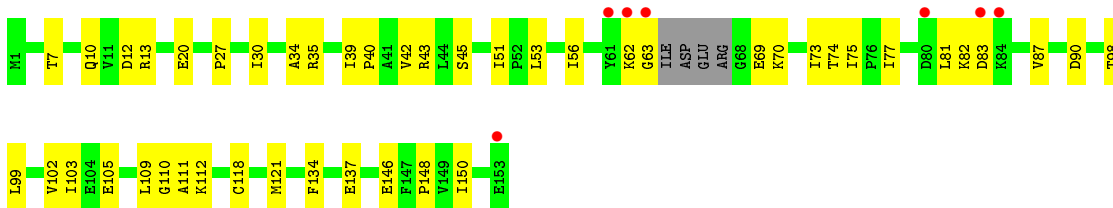
- Molecule 1: purine phosphoribosyltransferase



- Molecule 1: purine phosphoribosyltransferase



- Molecule 1: purine phosphoribosyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	105.60Å 157.35Å 170.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.55 – 2.50 39.55 – 2.49	Depositor EDS
% Data completeness (in resolution range)	100.0 (39.55-2.50) 99.7 (39.55-2.49)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 2.48Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.209 , 0.241 0.209 , 0.241	Depositor DCC
R_{free} test set	4967 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	42.0	Xtrriage
Anisotropy	0.454	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 57.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15181	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.67% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.44	0/1256	0.68	0/1701
1	B	0.44	0/1236	0.63	0/1675
1	C	0.39	0/1187	0.64	0/1612
1	D	0.54	2/1222 (0.2%)	0.63	0/1658
1	E	0.43	0/1240	0.65	0/1681
1	F	0.41	0/1242	0.64	0/1685
1	G	0.43	0/1263	0.67	0/1714
1	H	0.42	0/1243	0.64	0/1684
1	I	0.40	0/1213	0.63	0/1647
1	J	0.42	0/1224	0.62	0/1661
1	K	0.46	0/1218	0.67	0/1653
1	L	0.40	0/1236	0.59	0/1675
All	All	0.43	2/14780 (0.0%)	0.64	0/20046

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	9	TRP	NE1-CE2	8.76	1.49	1.37
1	D	8	TRP	NE1-CE2	8.73	1.49	1.37

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1227	0	1284	47	0
1	B	1207	0	1265	36	0
1	C	1159	0	1225	36	0
1	D	1193	0	1256	38	0
1	E	1211	0	1273	40	0
1	F	1213	0	1274	19	0
1	G	1233	0	1294	51	0
1	H	1214	0	1275	44	0
1	I	1184	0	1243	33	0
1	J	1195	0	1256	25	0
1	K	1189	0	1253	54	0
1	L	1207	0	1265	39	0
2	A	46	0	0	0	0
2	B	69	0	0	1	0
2	C	52	0	0	0	0
2	D	57	0	0	1	0
2	E	92	0	0	2	0
2	F	98	0	0	0	0
2	G	52	0	0	1	0
2	H	51	0	0	0	0
2	I	52	0	0	0	0
2	J	59	0	0	1	0
2	K	52	0	0	1	0
2	L	69	0	0	0	0
All	All	15181	0	15163	425	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (425) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:LYS:HB3	1:A:71:PRO:HD2	1.28	1.07
1:C:70:LYS:HB2	1:C:71:PRO:HD3	1.43	1.00
1:J:30:ILE:HG13	1:J:51:ILE:HD11	1.45	0.97
1:G:66:GLU:HG2	1:G:67:ARG:N	1.82	0.94
1:C:39:ILE:HB	1:C:40:PRO:HD3	1.49	0.92
1:H:62:LYS:HE3	1:H:62:LYS:H	1.34	0.91
1:B:83:ASP:HA	1:B:112:LYS:HB2	1.51	0.89
1:G:66:GLU:HG2	1:G:67:ARG:H	1.33	0.88
1:A:69:GLU:HG3	1:A:70:LYS:H	1.40	0.87
1:K:30:ILE:HG13	1:K:51:ILE:HD11	1.56	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:39:ILE:HB	1:K:40:PRO:HD3	1.58	0.86
1:F:30:ILE:HG13	1:F:51:ILE:HD11	1.57	0.86
1:A:70:LYS:HB3	1:A:71:PRO:CD	2.07	0.85
1:I:39:ILE:HB	1:I:40:PRO:HD3	1.61	0.83
1:J:20:GLU:HA	1:J:20:GLU:OE2	1.79	0.81
1:B:62:LYS:HG2	1:B:69:GLU:OE2	1.81	0.80
1:A:70:LYS:CB	1:A:71:PRO:HD2	2.08	0.79
1:C:70:LYS:HB2	1:C:71:PRO:CD	2.14	0.78
1:G:66:GLU:HG2	1:G:67:ARG:HG3	1.66	0.78
1:H:6:LEU:O	1:H:143:PRO:HG3	1.85	0.77
1:E:63:GLY:H	1:E:69:GLU:HA	1.51	0.75
1:C:83:ASP:HA	1:C:112:LYS:HB2	1.69	0.73
1:K:38:LEU:HD21	1:K:55:VAL:HG13	1.71	0.73
1:C:4:VAL:HG11	1:E:1:MET:CE	2.20	0.71
1:A:69:GLU:HG3	1:A:70:LYS:N	2.06	0.71
1:A:81:LEU:HD13	1:A:86:VAL:HG21	1.73	0.71
1:B:6:LEU:O	1:B:143:PRO:HG3	1.92	0.69
1:H:30:ILE:HG13	1:H:51:ILE:HD11	1.75	0.69
1:E:6:LEU:O	1:E:143:PRO:HG3	1.94	0.68
1:H:57:ASP:HB3	1:H:74:THR:HB	1.75	0.68
1:C:99:LEU:O	1:C:103:ILE:HG13	1.93	0.68
1:L:99:LEU:O	1:L:103:ILE:HG13	1.93	0.68
1:K:81:LEU:HD13	1:K:86:VAL:HG21	1.76	0.67
1:D:23:ARG:HG3	1:D:23:ARG:HH11	1.58	0.67
1:E:63:GLY:N	1:E:69:GLU:HA	2.08	0.67
1:B:103:ILE:HD11	1:B:116:ILE:HD11	1.77	0.67
1:B:120:ALA:HB2	1:B:134:PHE:HB2	1.77	0.67
1:G:120:ALA:HB2	1:G:134:PHE:HB2	1.75	0.67
1:L:30:ILE:HD12	1:L:45:SER:HA	1.76	0.67
1:H:31:ILE:HD13	1:H:54:LYS:HB2	1.75	0.67
1:G:1:MET:HE1	1:K:4:VAL:HG11	1.78	0.66
1:K:121:MET:HE2	1:K:135:ARG:HG3	1.78	0.66
1:K:43:ARG:O	1:K:47:ILE:HG13	1.95	0.66
1:J:1:MET:CE	1:L:134:PHE:HB3	2.25	0.65
1:K:37:GLY:O	1:K:40:PRO:HD2	1.96	0.65
1:L:74:THR:HG22	1:L:75:ILE:HG13	1.78	0.65
1:I:99:LEU:O	1:I:103:ILE:HG13	1.97	0.65
1:D:68:GLY:HA3	1:L:70:LYS:HG2	1.78	0.65
1:G:39:ILE:HB	1:G:40:PRO:HD3	1.78	0.64
1:H:103:ILE:HD11	1:H:116:ILE:HD11	1.80	0.64
1:L:30:ILE:HG13	1:L:51:ILE:HD11	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:85:ARG:HH11	1:E:85:ARG:HB2	1.63	0.64
1:E:88:ILE:HD13	1:E:102:VAL:HG12	1.78	0.64
1:A:39:ILE:HB	1:A:40:PRO:HD3	1.79	0.64
1:A:121:MET:CE	1:A:128:VAL:HG13	2.28	0.64
1:C:75:ILE:HD11	1:D:53:LEU:O	1.97	0.64
1:C:70:LYS:CB	1:C:71:PRO:HD3	2.25	0.63
1:G:85:ARG:HH11	1:G:85:ARG:HG3	1.64	0.62
1:H:152:LYS:O	1:H:153:GLU:HB2	1.98	0.62
1:J:1:MET:HE1	1:L:134:PHE:HB3	1.82	0.62
1:A:121:MET:HE3	1:A:128:VAL:HG13	1.82	0.61
1:D:1:MET:HE1	1:F:134:PHE:HB3	1.83	0.61
1:G:88:ILE:HD11	1:G:106:VAL:HG21	1.82	0.61
1:E:61:TYR:O	1:E:69:GLU:HA	2.01	0.60
1:J:80:ASP:OD2	1:J:82:LYS:HB2	2.01	0.60
1:H:118:CYS:O	1:H:132:TYR:HA	2.02	0.60
1:A:67:ARG:O	1:A:67:ARG:HG3	2.00	0.60
1:E:63:GLY:HA3	1:E:69:GLU:N	2.16	0.60
1:B:12:ASP:HB3	1:E:9:TRP:CH2	2.37	0.60
1:H:62:LYS:H	1:H:62:LYS:CE	2.11	0.60
1:I:121:MET:HE2	1:I:135:ARG:HG3	1.83	0.60
1:A:106:VAL:HG12	1:A:111:ALA:HB2	1.84	0.60
1:K:30:ILE:CG1	1:K:51:ILE:HD11	2.31	0.60
1:I:60:PHE:HB3	1:I:69:GLU:HG2	1.84	0.60
1:C:4:VAL:HG11	1:E:1:MET:HE1	1.82	0.59
1:E:48:LEU:O	1:E:51:ILE:HG23	2.01	0.59
1:I:103:ILE:HD11	1:I:116:ILE:HD11	1.83	0.59
1:A:33:VAL:O	1:A:37:GLY:HA3	2.03	0.59
1:C:70:LYS:CB	1:C:71:PRO:CD	2.81	0.59
1:A:69:GLU:CG	1:A:70:LYS:H	2.12	0.59
1:B:57:ASP:HB3	1:B:74:THR:HB	1.85	0.59
1:G:85:ARG:HG3	1:G:85:ARG:NH1	2.17	0.59
1:E:20:GLU:HA	1:E:20:GLU:OE1	2.02	0.59
1:C:83:ASP:CA	1:C:112:LYS:HB2	2.31	0.59
1:D:1:MET:CE	1:F:134:PHE:HB3	2.33	0.58
1:E:95:THR:OG1	1:E:97:LYS:HG2	2.02	0.58
1:J:103:ILE:HD11	1:J:116:ILE:HD11	1.85	0.58
1:H:61:TYR:O	1:H:70:LYS:HD2	2.03	0.58
1:B:31:ILE:HD13	1:B:54:LYS:HB2	1.86	0.58
1:C:39:ILE:HD13	1:D:39:ILE:HG12	1.85	0.58
1:A:94:ASP:OD2	1:A:122:LYS:HD2	2.04	0.57
1:E:122:LYS:HG2	1:E:136:THR:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:187:HOH:O	1:F:146:GLU:HG2	2.04	0.57
1:L:82:LYS:HA	1:L:110:GLY:O	2.04	0.57
1:L:62:LYS:HD3	1:L:69:GLU:HB2	1.86	0.57
1:E:43:ARG:O	1:E:47:ILE:HG13	2.04	0.57
1:A:34:ALA:HB1	1:A:35:ARG:HA	1.86	0.57
1:G:33:VAL:O	1:G:37:GLY:HA3	2.04	0.57
1:B:43:ARG:O	1:B:47:ILE:HG13	2.04	0.56
1:A:30:ILE:HG13	1:A:51:ILE:HD11	1.87	0.56
1:G:106:VAL:HG12	1:G:111:ALA:HB2	1.87	0.56
1:H:31:ILE:CD1	1:H:54:LYS:HB2	2.35	0.56
1:G:102:VAL:O	1:G:106:VAL:HG23	2.04	0.56
1:H:30:ILE:CG1	1:H:51:ILE:HD11	2.35	0.56
1:H:83:ASP:HA	1:H:112:LYS:HB2	1.86	0.56
1:A:57:ASP:HB3	1:A:74:THR:HB	1.87	0.56
1:B:99:LEU:O	1:B:103:ILE:HG13	2.05	0.56
1:B:80:ASP:OD2	1:B:82:LYS:HB2	2.06	0.56
1:D:68:GLY:CA	1:L:70:LYS:HG2	2.35	0.56
1:J:25:TYR:O	1:J:85:ARG:NH1	2.39	0.56
1:D:118:CYS:O	1:D:132:TYR:HA	2.06	0.56
1:K:11:VAL:O	1:K:15:ILE:HG13	2.06	0.55
1:B:3:LYS:HG2	1:B:139:TRP:HB3	1.88	0.55
1:K:120:ALA:HB2	1:K:134:PHE:HB2	1.89	0.54
1:D:23:ARG:HG3	1:D:23:ARG:NH1	2.22	0.54
1:A:106:VAL:HG12	1:A:111:ALA:CB	2.37	0.54
1:C:11:VAL:HG21	1:C:143:PRO:HG2	1.89	0.54
1:G:104:GLU:O	1:G:108:LYS:HG3	2.07	0.54
1:A:12:ASP:OD1	1:A:43:ARG:NH1	2.39	0.54
1:C:39:ILE:HB	1:C:40:PRO:CD	2.31	0.54
1:D:68:GLY:N	1:L:70:LYS:HG2	2.23	0.54
1:D:74:THR:HG21	2:D:177:HOH:O	2.06	0.54
1:D:74:THR:HG22	1:D:75:ILE:HG13	1.89	0.54
1:H:82:LYS:O	1:H:83:ASP:HB3	2.07	0.54
1:K:88:ILE:HD13	1:K:102:VAL:HG12	1.90	0.54
1:E:37:GLY:O	1:E:40:PRO:HD2	2.07	0.53
1:I:74:THR:HG22	1:I:75:ILE:HG13	1.90	0.53
1:H:122:LYS:HG2	1:H:136:THR:O	2.08	0.53
1:L:7:THR:OG1	1:L:10:GLN:HG3	2.09	0.53
1:H:30:ILE:CD1	1:H:51:ILE:HD11	2.38	0.53
1:G:37:GLY:O	1:G:40:PRO:HD2	2.09	0.53
1:A:99:LEU:O	1:A:103:ILE:HG13	2.09	0.53
1:H:51:ILE:HB	1:H:52:PRO:HD2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:83:ASP:O	1:D:112:LYS:HD2	2.08	0.53
1:K:91:ASP:OD1	1:K:92:VAL:N	2.40	0.53
1:E:83:ASP:CG	1:E:83:ASP:O	2.48	0.53
1:G:4:VAL:HG11	1:I:1:MET:CE	2.38	0.53
1:G:60:PHE:CE1	1:G:97:LYS:HE3	2.44	0.53
1:J:150:ILE:N	1:J:150:ILE:HD12	2.23	0.53
1:D:30:ILE:CD1	1:D:51:ILE:HD11	2.39	0.52
1:L:30:ILE:HG13	1:L:51:ILE:CD1	2.40	0.52
1:L:150:ILE:N	1:L:150:ILE:HD12	2.25	0.52
1:A:77:ILE:HG13	1:A:109:LEU:HD11	1.89	0.52
1:F:118:CYS:O	1:F:132:TYR:HA	2.09	0.52
1:G:59:LYS:NZ	1:G:74:THR:HG21	2.24	0.52
1:E:39:ILE:HB	1:E:40:PRO:HD3	1.91	0.52
1:I:39:ILE:CB	1:I:40:PRO:HD3	2.37	0.52
1:A:70:LYS:CB	1:A:71:PRO:CD	2.77	0.52
1:J:94:ASP:OD2	1:J:122:LYS:HD2	2.09	0.52
1:D:94:ASP:OD2	1:D:122:LYS:HD2	2.10	0.52
1:E:33:VAL:O	1:E:37:GLY:HA3	2.10	0.52
1:D:136:THR:HG22	1:D:140:ILE:HD11	1.92	0.52
1:G:92:VAL:HA	1:G:120:ALA:O	2.09	0.52
1:J:118:CYS:O	1:J:132:TYR:HA	2.09	0.52
1:I:8:TRP:CE2	1:I:143:PRO:HB2	2.45	0.52
1:G:43:ARG:O	1:G:47:ILE:HG13	2.10	0.51
1:I:30:ILE:HG13	1:I:51:ILE:HD11	1.92	0.51
1:F:39:ILE:HB	1:F:40:PRO:HD3	1.91	0.51
1:H:3:LYS:HG2	1:H:139:TRP:HB3	1.91	0.51
1:I:94:ASP:CG	1:I:122:LYS:HD2	2.30	0.51
1:G:6:LEU:O	1:G:143:PRO:HG3	2.10	0.51
1:I:39:ILE:HD12	1:I:39:ILE:N	2.25	0.51
1:B:30:ILE:HD11	1:B:51:ILE:HD11	1.92	0.51
1:J:30:ILE:CG1	1:J:51:ILE:HD11	2.29	0.51
1:B:92:VAL:HA	1:B:120:ALA:O	2.11	0.51
1:H:30:ILE:HD12	1:H:45:SER:HA	1.91	0.51
1:H:60:PHE:CE2	1:H:71:PRO:HB3	2.46	0.51
1:D:25:TYR:HD2	1:D:27:PRO:HG3	1.75	0.51
1:K:75:ILE:HD11	1:L:53:LEU:O	2.11	0.51
1:B:37:GLY:O	1:B:40:PRO:HD2	2.11	0.50
1:C:4:VAL:HG11	1:E:1:MET:HE3	1.93	0.50
1:D:91:ASP:OD2	1:D:92:VAL:N	2.44	0.50
1:I:60:PHE:CZ	1:I:97:LYS:HG2	2.46	0.50
1:H:43:ARG:O	1:H:47:ILE:HG13	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:39:ILE:HB	1:D:40:PRO:HD3	1.92	0.50
1:E:108:LYS:HG2	1:E:108:LYS:O	2.10	0.50
1:J:20:GLU:CA	1:J:20:GLU:OE2	2.53	0.50
1:K:77:ILE:O	1:K:77:ILE:HG13	2.12	0.50
1:G:118:CYS:O	1:G:132:TYR:HA	2.11	0.50
1:I:9:TRP:CE3	1:I:9:TRP:HA	2.47	0.50
1:J:57:ASP:HB3	1:J:74:THR:OG1	2.11	0.50
1:E:99:LEU:O	1:E:103:ILE:HG13	2.12	0.50
1:K:121:MET:HE3	1:K:122:LYS:C	2.31	0.50
1:F:25:TYR:HD2	1:F:27:PRO:HG3	1.76	0.50
1:H:70:LYS:HD2	1:H:70:LYS:H	1.77	0.50
1:A:30:ILE:HG13	1:A:51:ILE:CD1	2.42	0.49
1:A:6:LEU:O	1:A:143:PRO:HG3	2.12	0.49
1:C:102:VAL:O	1:C:106:VAL:HG23	2.12	0.49
1:E:63:GLY:H	1:E:69:GLU:CA	2.22	0.49
1:L:39:ILE:HB	1:L:40:PRO:HD3	1.93	0.49
1:G:59:LYS:HZ2	1:G:74:THR:HG21	1.76	0.49
1:E:74:THR:HG22	1:E:75:ILE:HG13	1.93	0.49
1:J:1:MET:HE2	1:L:134:PHE:HB3	1.94	0.49
1:A:106:VAL:CG1	1:A:111:ALA:HB2	2.42	0.49
1:G:8:TRP:CE2	1:G:143:PRO:HB2	2.47	0.49
1:J:99:LEU:O	1:J:103:ILE:HG13	2.13	0.49
1:G:79:GLY:O	1:G:109:LEU:HD22	2.13	0.49
1:K:6:LEU:O	1:K:143:PRO:HG3	2.13	0.49
1:K:83:ASP:HA	1:K:112:LYS:HB2	1.95	0.49
1:F:31:ILE:HD12	1:F:86:VAL:HG11	1.95	0.49
1:F:38:LEU:O	1:F:42:VAL:HG23	2.13	0.49
1:I:4:VAL:HG11	1:K:1:MET:HE1	1.95	0.49
1:D:68:GLY:HA3	1:L:70:LYS:CG	2.43	0.49
1:A:104:GLU:O	1:A:108:LYS:HB2	2.13	0.48
1:A:25:TYR:HD2	1:A:27:PRO:HG3	1.78	0.48
1:G:88:ILE:HD13	1:G:102:VAL:HG12	1.94	0.48
1:A:120:ALA:HB2	1:A:134:PHE:HB2	1.94	0.48
1:E:153:GLU:CD	1:E:153:GLU:H	2.16	0.48
1:J:39:ILE:HB	1:J:40:PRO:HD3	1.95	0.48
1:A:34:ALA:CB	1:A:35:ARG:HA	2.40	0.48
1:C:38:LEU:HD21	1:C:55:VAL:HG13	1.94	0.48
1:I:43:ARG:NH2	1:L:13:ARG:NH2	2.62	0.48
1:G:1:MET:CE	1:K:4:VAL:HG11	2.42	0.48
1:G:99:LEU:O	1:G:103:ILE:HG13	2.14	0.48
1:H:86:VAL:HG12	1:H:87:VAL:N	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:39:ILE:CD1	1:L:39:ILE:HG23	2.44	0.48
1:B:118:CYS:O	1:B:132:TYR:HA	2.13	0.48
1:D:103:ILE:HD11	1:D:116:ILE:HD11	1.95	0.48
1:K:85:ARG:HG3	1:K:85:ARG:NH1	2.29	0.48
1:L:81:LEU:O	1:L:111:ALA:HA	2.14	0.47
1:L:12:ASP:OD1	1:L:43:ARG:NH1	2.47	0.47
1:L:56:ILE:HA	1:L:75:ILE:O	2.14	0.47
1:G:150:ILE:HD12	1:K:132:TYR:CE1	2.49	0.47
1:I:4:VAL:HG11	1:K:1:MET:CE	2.44	0.47
1:K:2:ASP:O	1:K:138:LYS:HD3	2.13	0.47
1:L:77:ILE:HB	1:L:109:LEU:HD11	1.96	0.47
1:H:150:ILE:HD12	1:H:150:ILE:N	2.29	0.47
1:G:43:ARG:NH2	1:G:47:ILE:HD11	2.29	0.47
1:H:99:LEU:O	1:H:103:ILE:HG13	2.14	0.47
1:H:30:ILE:HG13	1:H:51:ILE:CD1	2.43	0.47
1:K:118:CYS:O	1:K:132:TYR:HA	2.15	0.47
1:G:39:ILE:CB	1:G:40:PRO:HD3	2.44	0.47
1:A:94:ASP:CG	1:A:122:LYS:HD2	2.35	0.47
1:K:60:PHE:CZ	1:K:97:LYS:HG2	2.49	0.47
1:L:62:LYS:HG3	1:L:63:GLY:H	1.79	0.47
1:C:9:TRP:HA	1:C:9:TRP:CE3	2.49	0.47
1:G:111:ALA:HB3	1:G:114:ILE:HD11	1.96	0.47
1:A:39:ILE:CB	1:A:40:PRO:HD3	2.44	0.47
1:D:30:ILE:HG13	1:D:51:ILE:HD11	1.96	0.47
1:E:118:CYS:O	1:E:132:TYR:HA	2.14	0.47
1:F:99:LEU:O	1:F:103:ILE:HG13	2.15	0.47
1:H:85:ARG:HH11	1:H:112:LYS:NZ	2.13	0.47
1:D:88:ILE:HD13	1:D:102:VAL:HG12	1.97	0.46
1:F:60:PHE:CE2	1:F:71:PRO:HB3	2.50	0.46
1:H:39:ILE:HB	1:H:40:PRO:HD3	1.97	0.46
1:I:39:ILE:HD12	1:I:39:ILE:H	1.80	0.46
1:A:30:ILE:CG1	1:A:51:ILE:HD11	2.46	0.46
1:C:25:TYR:O	1:C:26:LYS:HB2	2.16	0.46
1:I:85:ARG:HG3	1:I:85:ARG:HH11	1.80	0.46
1:A:122:LYS:HG2	1:A:136:THR:O	2.16	0.46
1:B:30:ILE:CD1	1:B:51:ILE:HD11	2.45	0.46
1:C:37:GLY:O	1:C:40:PRO:HD2	2.14	0.46
1:K:85:ARG:HH11	1:K:85:ARG:CG	2.28	0.46
1:G:83:ASP:CA	1:G:112:LYS:HB2	2.46	0.46
1:H:25:TYR:HD2	1:H:27:PRO:HG3	1.81	0.46
1:A:90:ASP:O	1:A:118:CYS:HA	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:83:ASP:HA	1:G:112:LYS:HB2	1.97	0.46
1:I:9:TRP:HE3	1:I:9:TRP:HA	1.80	0.46
1:L:42:VAL:O	1:L:45:SER:HB3	2.16	0.46
1:D:98:THR:O	1:D:102:VAL:HG23	2.16	0.45
1:A:69:GLU:CG	1:A:70:LYS:N	2.73	0.45
1:E:92:VAL:HG23	1:E:120:ALA:C	2.36	0.45
1:G:103:ILE:HD11	1:G:116:ILE:HD11	1.99	0.45
1:I:118:CYS:O	1:I:132:TYR:HA	2.15	0.45
1:D:138:LYS:HA	1:L:137:GLU:OE2	2.16	0.45
1:D:86:VAL:HG12	1:D:87:VAL:N	2.31	0.45
1:E:97:LYS:HG3	1:E:98:THR:N	2.32	0.45
1:F:86:VAL:HG12	1:F:87:VAL:N	2.31	0.45
1:G:106:VAL:HG12	1:G:111:ALA:CB	2.47	0.45
1:A:48:LEU:O	1:A:51:ILE:HG23	2.16	0.45
1:B:62:LYS:HA	1:B:69:GLU:HG2	1.99	0.45
1:D:102:VAL:O	1:D:106:VAL:HG23	2.16	0.45
1:G:57:ASP:HB3	1:G:74:THR:HB	1.98	0.45
1:L:73:ILE:HG22	1:L:73:ILE:O	2.16	0.45
1:E:121:MET:HG2	1:E:132:TYR:CE1	2.52	0.45
1:G:66:GLU:HG2	1:G:67:ARG:CG	2.44	0.45
1:H:12:ASP:HB3	1:K:9:TRP:CH2	2.52	0.45
1:D:81:LEU:HD13	1:D:86:VAL:HG21	1.98	0.45
1:J:82:LYS:O	1:J:83:ASP:HB2	2.17	0.45
1:L:34:ALA:HA	1:L:35:ARG:HA	1.59	0.45
1:C:88:ILE:HD13	1:C:102:VAL:HG12	1.99	0.45
1:H:30:ILE:HD11	1:H:51:ILE:HD11	1.97	0.45
1:I:4:VAL:CG1	1:K:1:MET:HE1	2.47	0.45
1:F:35:ARG:HD2	1:F:35:ARG:HH11	1.64	0.45
1:B:123:PRO:HG3	1:B:135:ARG:HE	1.81	0.44
1:B:30:ILE:HD12	1:B:45:SER:HA	1.99	0.44
1:D:101:VAL:HG12	1:D:101:VAL:O	2.16	0.44
1:D:25:TYR:CE2	1:D:87:VAL:HG22	2.52	0.44
1:K:8:TRP:CE2	1:K:143:PRO:HB2	2.51	0.44
1:C:87:VAL:O	1:C:87:VAL:HG13	2.16	0.44
1:H:34:ALA:HA	1:H:35:ARG:HA	1.62	0.44
1:H:37:GLY:O	1:H:40:PRO:HD2	2.16	0.44
1:I:29:VAL:HG23	1:I:52:PRO:O	2.17	0.44
1:G:83:ASP:HA	1:G:112:LYS:HD2	1.99	0.44
1:H:103:ILE:HD11	1:H:116:ILE:CD1	2.48	0.44
1:K:91:ASP:HB2	1:K:142:PHE:CZ	2.52	0.44
1:K:97:LYS:HB3	1:K:97:LYS:NZ	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:30:ILE:HG13	1:K:51:ILE:CD1	2.38	0.44
1:K:39:ILE:HB	1:K:40:PRO:CD	2.38	0.44
1:B:37:GLY:C	1:B:40:PRO:HD2	2.37	0.44
1:C:29:VAL:HG22	1:C:30:ILE:N	2.31	0.44
1:G:4:VAL:HG11	1:I:1:MET:HE3	2.00	0.44
1:B:1:MET:CE	1:D:134:PHE:HB3	2.47	0.44
1:B:60:PHE:CE2	1:B:71:PRO:HB3	2.52	0.44
1:J:30:ILE:HD12	1:J:45:SER:HA	1.99	0.44
2:G:656:HOH:O	1:H:78:HIS:HB3	2.18	0.44
1:K:73:ILE:CD1	1:K:105:GLU:HG3	2.47	0.44
1:F:30:ILE:HD12	1:F:45:SER:HA	2.00	0.44
1:A:84:LYS:O	1:A:112:LYS:HB3	2.18	0.44
1:B:146:GLU:HA	1:B:146:GLU:OE1	2.18	0.44
1:C:147:PHE:HA	1:C:148:PRO:HD3	1.83	0.43
1:H:62:LYS:N	1:H:62:LYS:CE	2.80	0.43
1:C:118:CYS:O	1:C:132:TYR:HA	2.18	0.43
1:D:27:PRO:HB3	1:D:87:VAL:HG23	2.00	0.43
1:E:8:TRP:CE2	1:E:143:PRO:HB2	2.53	0.43
1:F:27:PRO:HB3	1:F:87:VAL:HG23	1.98	0.43
1:I:121:MET:HG2	1:I:132:TYR:CD1	2.52	0.43
1:B:56:ILE:HB	1:B:77:ILE:HG23	2.00	0.43
1:G:91:ASP:OD1	1:G:91:ASP:N	2.48	0.43
1:K:25:TYR:CD2	1:K:27:PRO:HG3	2.53	0.43
1:A:118:CYS:O	1:A:132:TYR:HA	2.17	0.43
1:K:39:ILE:HG13	1:K:144:TRP:CD2	2.53	0.43
1:B:135:ARG:HD3	1:F:2:ASP:OD2	2.18	0.43
1:I:81:LEU:O	1:I:111:ALA:HA	2.19	0.43
1:E:121:MET:HG2	1:E:132:TYR:CD1	2.53	0.43
2:E:179:HOH:O	1:F:35:ARG:HD3	2.19	0.43
1:I:141:VAL:HG11	1:I:147:PHE:CE1	2.54	0.43
1:L:98:THR:O	1:L:102:VAL:HG23	2.18	0.43
1:B:83:ASP:CA	1:B:112:LYS:HB2	2.36	0.43
1:D:136:THR:CG2	1:D:140:ILE:HD11	2.48	0.43
1:G:90:ASP:O	1:G:118:CYS:HA	2.19	0.43
1:H:152:LYS:O	1:H:153:GLU:CB	2.65	0.43
1:K:136:THR:HG21	1:K:140:ILE:HD11	2.00	0.43
1:G:1:MET:HE3	1:K:4:VAL:HG21	2.00	0.43
1:C:39:ILE:HG13	1:C:144:TRP:CD2	2.53	0.43
1:E:120:ALA:HB2	1:E:134:PHE:HB2	2.00	0.43
1:J:115:LYS:HG2	1:J:131:TYR:HE1	1.83	0.43
1:K:87:VAL:O	1:K:87:VAL:HG13	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:88:ILE:HD11	1:K:106:VAL:HG21	2.01	0.43
1:L:146:GLU:O	1:L:148:PRO:HD3	2.19	0.43
1:C:150:ILE:HD12	1:C:150:ILE:N	2.34	0.43
1:C:9:TRP:HA	1:C:9:TRP:HE3	1.84	0.43
1:E:86:VAL:HG12	1:E:87:VAL:N	2.34	0.43
1:K:25:TYR:HD2	1:K:27:PRO:HG3	1.84	0.43
1:A:59:LYS:HE3	1:A:59:LYS:HB2	1.80	0.42
1:B:121:MET:HG2	1:B:132:TYR:CE2	2.54	0.42
1:J:86:VAL:HG12	1:J:87:VAL:N	2.33	0.42
1:L:73:ILE:HD13	1:L:105:GLU:HG2	2.00	0.42
1:A:8:TRP:CE2	1:A:143:PRO:HB2	2.54	0.42
1:G:122:LYS:HG2	1:G:136:THR:O	2.18	0.42
1:H:153:GLU:CA	1:H:153:GLU:OE1	2.66	0.42
1:H:62:LYS:HE3	1:H:62:LYS:N	2.16	0.42
1:K:43:ARG:NH2	2:K:171:HOH:O	2.41	0.42
1:K:29:VAL:CG1	1:K:86:VAL:HG22	2.49	0.42
1:L:83:ASP:C	1:L:112:LYS:HB2	2.39	0.42
1:E:60:PHE:CE2	1:E:71:PRO:HB3	2.55	0.42
1:K:136:THR:CG2	1:K:140:ILE:HD11	2.50	0.42
1:B:121:MET:HG2	1:B:132:TYR:CD2	2.55	0.42
1:C:38:LEU:HD12	1:D:42:VAL:CG2	2.50	0.42
1:J:103:ILE:HD11	1:J:116:ILE:CD1	2.50	0.42
1:L:90:ASP:O	1:L:118:CYS:HA	2.19	0.42
1:A:62:LYS:HD3	1:A:63:GLY:H	1.84	0.42
1:H:100:GLU:HG3	1:H:127:VAL:CG2	2.50	0.42
1:K:121:MET:CE	1:K:135:ARG:HG3	2.47	0.42
1:B:97:LYS:HG3	1:B:98:THR:N	2.35	0.42
1:B:33:VAL:HG21	1:B:102:VAL:HG21	2.02	0.42
1:I:39:ILE:CD1	1:I:39:ILE:H	2.33	0.42
1:L:27:PRO:HB3	1:L:87:VAL:HG23	2.00	0.42
1:B:94:ASP:OD2	1:B:122:LYS:HD2	2.19	0.42
1:H:92:VAL:HA	1:H:120:ALA:O	2.20	0.42
1:K:39:ILE:HD12	1:L:39:ILE:HG23	2.02	0.42
1:C:141:VAL:HG11	1:C:147:PHE:CE1	2.54	0.42
1:C:141:VAL:HG11	1:C:147:PHE:HE1	1.84	0.42
1:E:12:ASP:OD1	1:E:43:ARG:NH1	2.50	0.42
1:E:88:ILE:HD13	1:E:102:VAL:CG1	2.48	0.42
1:H:153:GLU:HA	1:H:153:GLU:OE1	2.19	0.42
1:K:99:LEU:HD23	1:K:99:LEU:HA	1.84	0.42
1:B:85:ARG:NH1	2:B:205:HOH:O	2.51	0.42
1:B:135:ARG:NH1	1:J:135:ARG:NH1	2.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:ASP:HB2	1:A:75:ILE:HD12	2.01	0.41
1:G:6:LEU:HA	1:G:6:LEU:HD23	1.90	0.41
1:G:63:GLY:HA3	1:G:70:LYS:CE	2.49	0.41
1:D:25:TYR:HE2	1:D:87:VAL:HG22	1.85	0.41
1:I:22:LEU:O	1:I:27:PRO:HD3	2.20	0.41
1:K:121:MET:HE3	1:K:123:PRO:N	2.35	0.41
1:D:27:PRO:HB3	1:D:87:VAL:CG2	2.50	0.41
1:E:83:ASP:OD1	1:E:83:ASP:O	2.38	0.41
1:H:9:TRP:CZ3	1:L:7:THR:HA	2.55	0.41
1:K:30:ILE:HG12	1:K:87:VAL:CG1	2.51	0.41
1:L:30:ILE:CG1	1:L:51:ILE:HD11	2.50	0.41
1:A:115:LYS:HG2	1:A:131:TYR:HE1	1.86	0.41
1:B:34:ALA:HA	1:B:35:ARG:HA	1.78	0.41
1:G:39:ILE:HG22	1:G:40:PRO:N	2.36	0.41
1:I:97:LYS:HB3	1:I:97:LYS:HE2	1.88	0.41
1:J:67:ARG:NH2	2:J:154:HOH:O	2.53	0.41
1:K:81:LEU:HA	1:K:84:LYS:HD2	2.01	0.41
1:C:25:TYR:HD2	1:C:27:PRO:HG3	1.85	0.41
1:E:34:ALA:HA	1:E:35:ARG:HA	1.63	0.41
1:G:135:ARG:HH11	1:G:135:ARG:HD3	1.60	0.41
1:A:60:PHE:CE2	1:A:71:PRO:HB3	2.56	0.41
1:I:34:ALA:HA	1:I:35:ARG:HA	1.59	0.41
1:L:27:PRO:HB3	1:L:87:VAL:CG2	2.50	0.41
1:A:22:LEU:O	1:A:27:PRO:HD3	2.20	0.41
1:B:81:LEU:HD13	1:B:86:VAL:HG21	2.02	0.41
1:D:9:TRP:HZ2	1:D:13:ARG:NH1	2.19	0.41
1:E:99:LEU:HA	1:E:99:LEU:HD23	1.86	0.41
1:A:88:ILE:HD11	1:A:106:VAL:HG21	2.03	0.41
1:A:6:LEU:HA	1:A:6:LEU:HD23	1.87	0.41
1:C:38:LEU:O	1:C:42:VAL:HG23	2.20	0.41
1:C:56:ILE:O	1:C:56:ILE:HG23	2.21	0.41
1:G:26:LYS:N	1:G:27:PRO:HD3	2.36	0.41
1:G:59:LYS:HZ2	1:G:74:THR:CG2	2.34	0.41
1:K:85:ARG:NH1	1:K:85:ARG:CG	2.84	0.41
1:K:9:TRP:HA	1:K:9:TRP:CE3	2.56	0.41
1:E:94:ASP:OD2	1:E:122:LYS:HD2	2.20	0.41
1:G:103:ILE:O	1:G:107:LYS:HG3	2.21	0.40
1:G:30:ILE:HD11	1:G:51:ILE:HD11	2.03	0.40
1:H:86:VAL:CG1	1:H:87:VAL:N	2.84	0.40
1:C:42:VAL:O	1:C:45:SER:HB3	2.21	0.40
1:F:138:LYS:HA	1:J:137:GLU:OE2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:75:ILE:HD11	1:J:53:LEU:O	2.22	0.40
1:F:30:ILE:HG13	1:F:51:ILE:CD1	2.38	0.40
1:I:58:VAL:HB	1:I:98:THR:HG23	2.03	0.40
1:C:25:TYR:CD2	1:C:27:PRO:HG3	2.57	0.40
1:D:150:ILE:HD12	1:D:150:ILE:N	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	147/153 (96%)	135 (92%)	10 (7%)	2 (1%)	11	20
1	B	145/153 (95%)	139 (96%)	6 (4%)	0	100	100
1	C	139/153 (91%)	129 (93%)	10 (7%)	0	100	100
1	D	143/153 (94%)	133 (93%)	8 (6%)	2 (1%)	11	20
1	E	145/153 (95%)	138 (95%)	7 (5%)	0	100	100
1	F	146/153 (95%)	143 (98%)	3 (2%)	0	100	100
1	G	150/153 (98%)	144 (96%)	5 (3%)	1 (1%)	22	39
1	H	145/153 (95%)	139 (96%)	6 (4%)	0	100	100
1	I	142/153 (93%)	135 (95%)	7 (5%)	0	100	100
1	J	143/153 (94%)	138 (96%)	5 (4%)	0	100	100
1	K	142/153 (93%)	135 (95%)	7 (5%)	0	100	100
1	L	145/153 (95%)	140 (97%)	5 (3%)	0	100	100
All	All	1732/1836 (94%)	1648 (95%)	79 (5%)	5 (0%)	41	61

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	152	LYS
1	G	112	LYS
1	A	70	LYS
1	D	110	GLY
1	D	73	ILE

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	134/136 (98%)	128 (96%)	6 (4%)	27	51
1	B	132/136 (97%)	130 (98%)	2 (2%)	65	85
1	C	128/136 (94%)	126 (98%)	2 (2%)	62	84
1	D	131/136 (96%)	129 (98%)	2 (2%)	65	85
1	E	133/136 (98%)	129 (97%)	4 (3%)	41	68
1	F	133/136 (98%)	131 (98%)	2 (2%)	65	85
1	G	135/136 (99%)	130 (96%)	5 (4%)	34	60
1	H	133/136 (98%)	128 (96%)	5 (4%)	33	58
1	I	130/136 (96%)	127 (98%)	3 (2%)	50	76
1	J	131/136 (96%)	125 (95%)	6 (5%)	27	50
1	K	131/136 (96%)	128 (98%)	3 (2%)	50	76
1	L	132/136 (97%)	130 (98%)	2 (2%)	65	85
All	All	1583/1632 (97%)	1541 (97%)	42 (3%)	44	71

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASP
1	A	35	ARG
1	A	62	LYS
1	A	67	ARG
1	A	70	LYS
1	A	137	GLU

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Mol	Chain	Res	Type
1	B	16	PHE
1	B	62	LYS
1	C	12	ASP
1	C	121	MET
1	D	20	GLU
1	D	50	ASP
1	E	59	LYS
1	E	62	LYS
1	E	85	ARG
1	E	108	LYS
1	F	35	ARG
1	F	121	MET
1	G	35	ARG
1	G	59	LYS
1	G	65	ASP
1	G	74	THR
1	G	121	MET
1	H	62	LYS
1	H	70	LYS
1	H	80	ASP
1	H	121	MET
1	H	153	GLU
1	I	35	ARG
1	I	104	GLU
1	I	121	MET
1	J	16	PHE
1	J	67	ARG
1	J	80	ASP
1	J	113	GLU
1	J	115	LYS
1	J	121	MET
1	K	6	LEU
1	K	85	ARG
1	K	121	MET
1	L	20	GLU
1	L	121	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data i

6.1 Protein, DNA and RNA chains i

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	151/153 (98%)	0.12	10 (6%) 18 19	24, 46, 75, 104	0
1	B	149/153 (97%)	0.08	5 (3%) 45 48	23, 41, 69, 91	0
1	C	143/153 (93%)	0.25	7 (4%) 29 31	24, 51, 81, 85	0
1	D	147/153 (96%)	0.26	9 (6%) 21 22	25, 48, 87, 97	0
1	E	149/153 (97%)	0.05	7 (4%) 31 33	23, 40, 68, 88	0
1	F	150/153 (98%)	-0.13	0 100 100	22, 38, 59, 82	0
1	G	152/153 (99%)	0.13	3 (1%) 65 68	23, 48, 78, 89	0
1	H	149/153 (97%)	0.05	5 (3%) 45 48	23, 44, 80, 86	0
1	I	146/153 (95%)	-0.06	2 (1%) 75 77	20, 42, 64, 79	0
1	J	147/153 (96%)	-0.20	2 (1%) 75 77	22, 37, 62, 75	0
1	K	146/153 (95%)	0.08	7 (4%) 30 32	23, 40, 73, 84	0
1	L	149/153 (97%)	0.10	7 (4%) 31 33	23, 46, 79, 89	0
All	All	1778/1836 (96%)	0.06	64 (3%) 42 46	20, 44, 78, 104	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	78	HIS	4.5
1	E	153	GLU	4.1
1	I	68	GLY	4.1
1	A	66	GLU	4.0
1	B	153	GLU	4.0
1	D	61	TYR	3.9
1	E	64	ILE	3.8
1	A	69	GLU	3.5
1	E	63	GLY	3.4
1	E	62	LYS	3.3
1	G	67	ARG	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	67	ARG	3.2
1	A	153	GLU	3.2
1	C	71	PRO	3.1
1	D	81	LEU	3.1
1	B	61	TYR	3.1
1	H	61	TYR	3.0
1	K	61	TYR	3.0
1	K	80	ASP	3.0
1	L	153	GLU	3.0
1	A	68	GLY	3.0
1	E	61	TYR	3.0
1	H	70	LYS	2.9
1	L	84	LYS	2.9
1	C	97	LYS	2.9
1	D	68	GLY	2.9
1	L	80	ASP	2.9
1	J	61	TYR	2.8
1	G	83	ASP	2.7
1	G	82	LYS	2.7
1	B	68	GLY	2.7
1	J	79	GLY	2.7
1	K	83	ASP	2.7
1	A	112	LYS	2.7
1	K	70	LYS	2.7
1	B	112	LYS	2.6
1	D	70	LYS	2.6
1	H	78	HIS	2.6
1	L	63	GLY	2.6
1	L	61	TYR	2.6
1	C	83	ASP	2.5
1	H	111	ALA	2.5
1	L	83	ASP	2.5
1	A	63	GLY	2.4
1	H	153	GLU	2.4
1	C	80	ASP	2.4
1	C	72	VAL	2.4
1	L	62	LYS	2.3
1	D	24	GLU	2.3
1	A	111	ALA	2.3
1	D	112	LYS	2.3
1	D	83	ASP	2.3
1	C	82	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	101	VAL	2.2
1	E	2	ASP	2.2
1	C	81	LEU	2.2
1	E	83	ASP	2.2
1	K	69	GLU	2.2
1	B	62	LYS	2.1
1	D	80	ASP	2.1
1	K	79	GLY	2.1
1	A	83	ASP	2.1
1	I	83	ASP	2.1
1	A	70	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.