



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:50 PM GMT

PDB ID : 1VG4
Title : Crystal Structure Of Octaprenyl Pyrophosphate Synthase From Hyperthermophilic Thermotoga Maritima F132A/L128A mutant
Authors : Guo, R.T.; Kuo, C.J.; Ko, T.P.; Chou, C.C.; Liang, P.H.; Wang, A.H.-J.
Deposited on : 2004-04-23
Resolution : 3.30 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

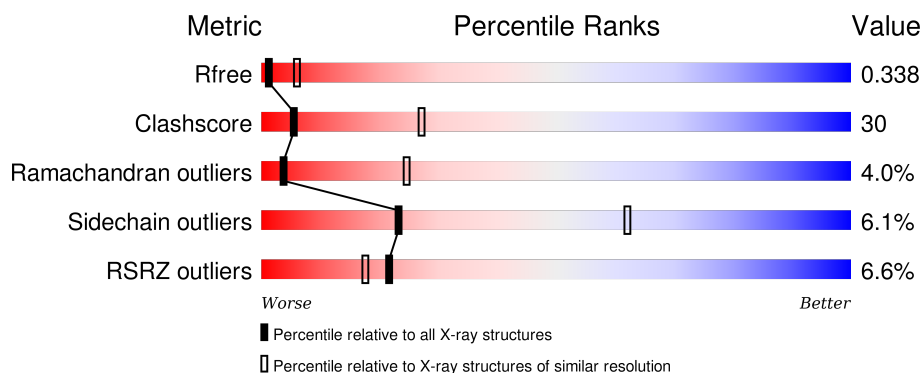
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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}	91344	2060 (3.40-3.20)
Clashscore	102246	1058 (3.38-3.22)
Ramachandran outliers	100387	1038 (3.38-3.22)
Sidechain outliers	100360	1037 (3.38-3.22)
RSRZ outliers	91569	2070 (3.40-3.20)

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. 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	299	
1	B	299	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4816 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 octoprenyl-diphosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	280	Total	C	N	O	S	0	0	0
			2219	1421	371	415	12			
1	B	280	Total	C	N	O	S	0	0	0
			2219	1421	371	415	12			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	128	ALA	LEU	ENGINEERED	UNP Q9X1M1
A	132	ALA	PHE	ENGINEERED	UNP Q9X1M1
B	128	ALA	LEU	ENGINEERED	UNP Q9X1M1
B	132	ALA	PHE	ENGINEERED	UNP Q9X1M1

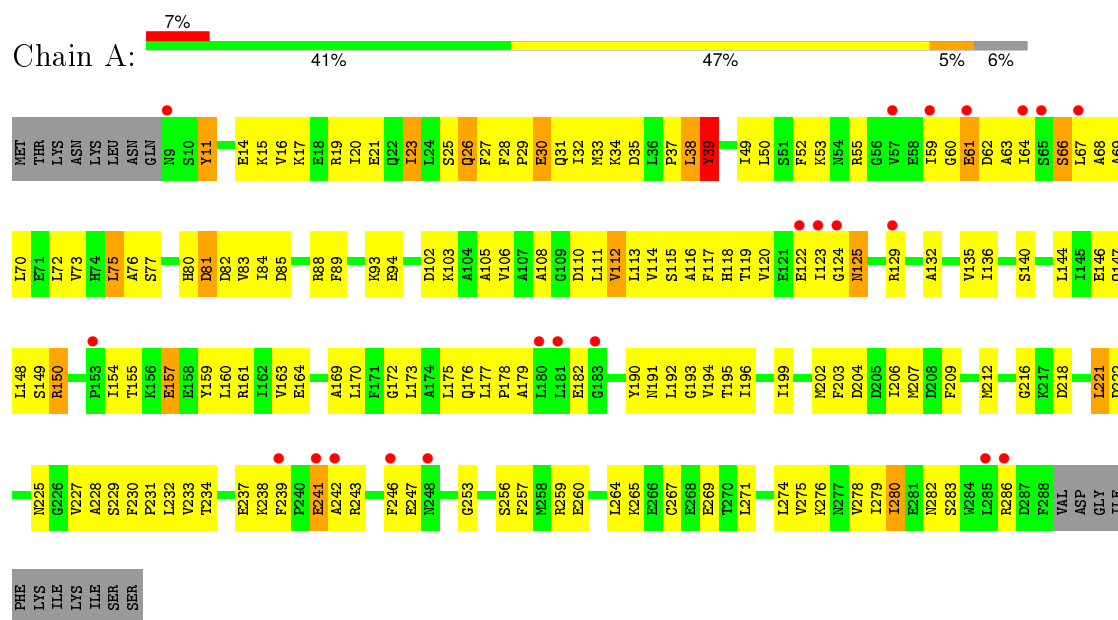
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	194	Total	O	0	0
			194	194		
2	B	184	Total	O	0	0
			184	184		

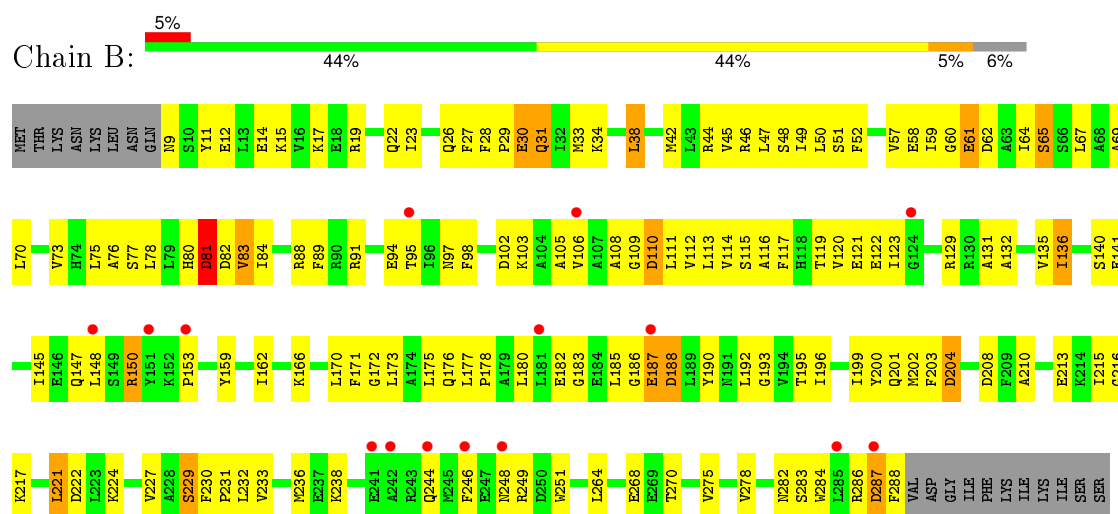
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: octoprenyl-diphosphate synthase



- Molecule 1: octoprenyl-diphosphate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	151.43 Å 151.43 Å 67.04 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.89 – 3.30 47.89 – 3.30	Depositor EDS
% Data completeness (in resolution range)	73.7 (47.89-3.30) 90.4 (47.89-3.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.16 (at 3.33 Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.277 , 0.321 0.295 , 0.338	Depositor DCC
R_{free} test set	578 reflections (5.23%)	DCC
Wilson B-factor (Å ²)	73.0	Xtriage
Anisotropy	0.569	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 80.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	1 of 11640 reflections (0.009%)	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	4816	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.24% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.28	0/2253	0.60	0/3025
1	B	0.28	0/2253	0.58	0/3025
All	All	0.28	0/4506	0.59	0/6050

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2219	0	2259	147	0
1	B	2219	0	2259	130	0
2	A	194	0	0	0	2
2	B	184	0	0	1	0
All	All	4816	0	4518	266	2

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 30.

All (266) 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:B:70:LEU:HD11	1:B:170:LEU:HD11	1.40	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:44:ARG:HE	1:B:170:LEU:HD23	1.26	0.97
1:B:244:GLN:HG2	1:B:248:ASN:HD21	1.29	0.96
1:A:15:LYS:HG2	1:A:64:ILE:HG21	1.52	0.91
1:A:77:SER:O	1:A:81:ASP:HB2	1.73	0.87
1:B:217:LYS:H	1:B:217:LYS:HD2	1.40	0.84
1:B:120:VAL:HG12	1:B:129:ARG:HG2	1.59	0.84
1:A:120:VAL:HG12	1:A:129:ARG:HG3	1.61	0.83
1:B:244:GLN:HG2	1:B:248:ASN:ND2	1.96	0.80
1:A:118:HIS:O	1:A:122:GLU:HG2	1.82	0.79
1:A:73:VAL:O	1:A:76:ALA:HB3	1.82	0.79
1:B:70:LEU:HD11	1:B:170:LEU:CD1	2.12	0.79
1:B:9:ASN:N	1:B:49:ILE:HG21	1.99	0.77
1:B:116:ALA:O	1:B:120:VAL:HG23	1.88	0.74
1:B:19:ARG:HG2	1:B:64:ILE:HG22	1.71	0.73
1:A:111:LEU:HD13	1:B:140:SER:HB2	1.69	0.72
1:A:276:LYS:O	1:A:280:ILE:HG12	1.89	0.72
1:B:102:ASP:O	1:B:106:VAL:HG23	1.90	0.71
1:B:217:LYS:HD2	1:B:217:LYS:N	2.06	0.71
1:A:175:LEU:HD11	1:A:196:ILE:HD12	1.73	0.69
1:A:238:LYS:HE2	1:A:239:PHE:HE1	1.56	0.69
1:A:216:GLY:H	1:A:222:ASP:HB2	1.58	0.68
1:A:160:LEU:O	1:A:164:GLU:HG3	1.94	0.68
1:A:238:LYS:HE2	1:A:239:PHE:CE1	2.28	0.68
1:A:75:LEU:HD12	1:A:75:LEU:H	1.58	0.68
1:A:88:ARG:O	1:A:94:GLU:HA	1.94	0.68
1:B:44:ARG:NE	1:B:170:LEU:HD23	2.06	0.67
1:B:12:GLU:HG2	1:B:59:ILE:HD12	1.76	0.67
1:A:110:ASP:O	1:A:114:VAL:HG23	1.95	0.66
1:A:135:VAL:HG21	1:A:173:LEU:HD22	1.75	0.66
1:A:216:GLY:N	1:A:222:ASP:HB2	2.11	0.66
1:A:116:ALA:O	1:A:120:VAL:HG23	1.96	0.66
1:A:62:ASP:O	1:A:66:SER:HB2	1.96	0.66
1:A:108:ALA:O	1:A:112:VAL:HG23	1.94	0.66
1:A:172:GLY:HA2	1:A:193:GLY:HA3	1.79	0.65
1:B:147:GLN:O	1:B:150:ARG:HB2	1.97	0.65
1:B:120:VAL:CG1	1:B:129:ARG:HG2	2.27	0.64
1:B:227:VAL:HG22	1:B:246:PHE:HZ	1.62	0.64
1:B:47:LEU:HD12	1:B:171:PHE:CD1	2.34	0.63
1:A:271:LEU:O	1:A:275:VAL:HG23	1.98	0.63
1:B:208:ASP:OD1	1:B:224:LYS:HG3	1.99	0.63
1:A:192:LEU:O	1:A:196:ILE:HG13	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:230:PHE:HB3	1:B:231:PRO:HD3	1.81	0.62
1:A:19:ARG:O	1:A:23:ILE:HG22	2.00	0.62
1:B:175:LEU:HD11	1:B:196:ILE:HD12	1.81	0.62
1:B:49:ILE:HG22	1:B:50:LEU:N	2.14	0.62
1:A:243:ARG:O	1:A:247:GLU:HG3	2.00	0.62
1:B:11:TYR:O	1:B:15:LYS:HG3	2.00	0.62
1:B:108:ALA:O	1:B:112:VAL:HG23	1.99	0.62
1:A:61:GLU:C	1:A:63:ALA:H	2.03	0.61
1:A:160:LEU:HA	1:A:163:VAL:HG22	1.81	0.61
1:B:69:ALA:HB2	1:B:119:THR:HG21	1.82	0.61
1:A:84:ILE:HD13	1:B:103:LYS:HG3	1.81	0.61
1:A:177:LEU:HB2	1:A:178:PRO:HD3	1.81	0.61
1:A:29:PRO:HG2	1:B:148:LEU:HD11	1.83	0.61
1:B:70:LEU:HD22	1:B:173:LEU:HD23	1.83	0.61
1:B:217:LYS:H	1:B:217:LYS:CD	2.14	0.60
1:B:175:LEU:O	1:B:178:PRO:HD2	2.02	0.60
1:B:182:GLU:HG3	2:B:385:HOH:O	2.01	0.60
1:A:69:ALA:HB2	1:A:119:THR:CG2	2.31	0.60
1:A:15:LYS:HG2	1:A:64:ILE:CG2	2.29	0.60
1:A:23:ILE:HD11	1:A:115:SER:O	2.02	0.60
1:B:12:GLU:O	1:B:15:LYS:HB2	2.02	0.60
1:B:113:LEU:O	1:B:116:ALA:HB3	2.02	0.59
1:A:111:LEU:HD13	1:B:140:SER:CB	2.33	0.59
1:B:131:ALA:O	1:B:135:VAL:HG23	2.02	0.59
1:A:21:GLU:HG3	1:A:33:MET:HE1	1.83	0.59
1:B:30:GLU:O	1:B:33:MET:HB3	2.02	0.59
1:A:227:VAL:C	1:A:229:SER:H	2.06	0.59
1:B:186:GLY:O	1:B:188:ASP:N	2.35	0.59
1:A:62:ASP:OD2	1:A:62:ASP:O	2.20	0.59
1:A:69:ALA:O	1:A:72:LEU:HB2	2.03	0.59
1:B:69:ALA:HB2	1:B:119:THR:CG2	2.33	0.58
1:B:264:LEU:O	1:B:268:GLU:HG3	2.03	0.58
1:A:132:ALA:O	1:A:136:ILE:HG13	2.04	0.58
1:A:27:PHE:HZ	1:A:115:SER:HA	1.68	0.58
1:B:69:ALA:O	1:B:73:VAL:HG23	2.03	0.58
1:A:106:VAL:HG21	1:B:84:ILE:HD11	1.85	0.58
1:B:47:LEU:HD12	1:B:171:PHE:HD1	1.69	0.57
1:A:20:ILE:HG12	1:A:68:ALA:HB1	1.85	0.57
1:B:97:ASN:HA	1:B:105:ALA:HB2	1.86	0.57
1:B:49:ILE:O	1:B:52:PHE:HB2	2.04	0.57
1:A:112:VAL:O	1:A:115:SER:N	2.28	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:119:THR:O	1:B:122:GLU:HB2	2.06	0.56
1:A:61:GLU:C	1:A:63:ALA:N	2.58	0.56
1:B:42:MET:O	1:B:46:ARG:HG3	2.05	0.56
1:B:110:ASP:O	1:B:114:VAL:HG23	2.06	0.56
1:A:11:TYR:C	1:A:11:TYR:CD1	2.78	0.56
1:A:113:LEU:HD23	1:B:113:LEU:HD23	1.86	0.56
1:B:227:VAL:HG22	1:B:246:PHE:CZ	2.41	0.56
1:B:244:GLN:CG	1:B:248:ASN:HD21	2.08	0.56
1:A:275:VAL:O	1:A:279:ILE:HG12	2.06	0.55
1:A:154:ILE:HD11	1:A:228:ALA:HA	1.89	0.55
1:A:69:ALA:HB2	1:A:119:THR:HG21	1.86	0.55
1:A:80:HIS:HE1	1:B:110:ASP:OD1	1.90	0.55
1:B:159:TYR:O	1:B:162:ILE:HB	2.06	0.55
1:A:253:GLY:O	1:A:256:SER:HB3	2.07	0.55
1:A:242:ALA:HB2	1:A:257:PHE:CD2	2.42	0.55
1:A:70:LEU:HD23	1:A:170:LEU:CD1	2.38	0.54
1:B:110:ASP:O	1:B:114:VAL:N	2.35	0.54
1:A:60:GLY:O	1:A:62:ASP:N	2.41	0.54
1:B:270:THR:HG22	1:B:270:THR:O	2.06	0.54
1:A:16:VAL:O	1:A:20:ILE:HG13	2.08	0.53
1:A:117:PHE:C	1:A:119:THR:N	2.62	0.53
1:B:176:GLN:HG2	1:B:180:LEU:HG	1.90	0.53
1:B:141:GLU:HG2	1:B:145:ILE:CD1	2.39	0.53
1:A:195:THR:O	1:A:199:ILE:HG13	2.09	0.53
1:B:88:ARG:HG3	1:B:98:PHE:CG	2.43	0.53
1:A:63:ALA:O	1:A:67:LEU:HG	2.09	0.53
1:B:76:ALA:HB1	1:B:109:GLY:O	2.08	0.53
1:A:23:ILE:O	1:A:26:GLN:HB2	2.09	0.53
1:A:62:ASP:CG	1:A:123:ILE:HG23	2.30	0.53
1:B:65:SER:HB3	1:B:123:ILE:HD11	1.90	0.52
1:A:206:ILE:HG23	1:A:264:LEU:HD11	1.90	0.52
1:A:120:VAL:HG12	1:A:120:VAL:O	2.10	0.52
1:B:178:PRO:O	1:B:182:GLU:HG2	2.09	0.52
1:B:51:SER:HA	1:B:284:TRP:CZ3	2.45	0.52
1:B:192:LEU:O	1:B:196:ILE:HG13	2.10	0.52
1:A:227:VAL:O	1:A:229:SER:N	2.37	0.51
1:A:140:SER:CB	1:B:111:LEU:HD13	2.40	0.51
1:A:140:SER:HB2	1:B:111:LEU:HD13	1.93	0.51
1:A:238:LYS:HB3	1:A:239:PHE:CD1	2.46	0.51
1:A:229:SER:HB3	1:A:232:LEU:CB	2.40	0.51
1:A:147:GLN:C	1:A:149:SER:H	2.13	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:SER:O	1:A:233:VAL:HG23	2.11	0.51
1:B:12:GLU:HG2	1:B:59:ILE:CD1	2.40	0.50
1:A:170:LEU:HD12	1:A:170:LEU:O	2.12	0.50
1:A:155:THR:OG1	1:A:157:GLU:HG3	2.11	0.50
1:B:185:LEU:HD13	1:B:187:GLU:OE1	2.13	0.49
1:A:178:PRO:O	1:A:182:GLU:HB2	2.12	0.49
1:A:192:LEU:HG	1:A:196:ILE:HD11	1.93	0.49
1:B:286:ARG:O	1:B:287:ASP:C	2.50	0.49
1:B:166:LYS:HE3	1:B:201:GLN:NE2	2.27	0.49
1:A:62:ASP:OD1	1:A:123:ILE:HG23	2.13	0.49
1:B:44:ARG:HH11	1:B:44:ARG:HG3	1.78	0.49
1:A:278:VAL:O	1:A:282:ASN:HB2	2.12	0.49
1:A:19:ARG:NH1	1:A:23:ILE:HG22	2.27	0.49
1:A:242:ALA:HB2	1:A:257:PHE:CE2	2.47	0.49
1:B:141:GLU:HG2	1:B:145:ILE:HD11	1.93	0.49
1:A:190:TYR:O	1:A:194:VAL:HG23	2.13	0.49
1:B:60:GLY:O	1:B:61:GLU:C	2.51	0.49
1:A:117:PHE:C	1:A:119:THR:H	2.16	0.48
1:B:221:LEU:O	1:B:222:ASP:HB3	2.12	0.48
1:A:227:VAL:C	1:A:229:SER:N	2.67	0.48
1:A:256:SER:O	1:A:260:GLU:HB2	2.13	0.48
1:A:117:PHE:O	1:A:120:VAL:N	2.47	0.48
1:A:89:PHE:HA	1:A:93:LYS:O	2.13	0.48
1:B:65:SER:HB3	1:B:123:ILE:CD1	2.44	0.48
1:B:77:SER:O	1:B:81:ASP:HB2	2.13	0.48
1:B:80:HIS:O	1:B:83:VAL:HG23	2.14	0.48
1:B:89:PHE:CE2	1:B:94:GLU:HG2	2.49	0.47
1:B:249:ARG:NE	1:B:251:TRP:CH2	2.79	0.47
1:B:19:ARG:O	1:B:23:ILE:HG13	2.15	0.47
1:B:132:ALA:O	1:B:136:ILE:HG12	2.14	0.47
1:B:57:VAL:HG12	1:B:58:GLU:N	2.30	0.47
1:A:49:ILE:O	1:A:53:LYS:HG2	2.15	0.47
1:B:45:VAL:HG13	1:B:67:LEU:HD13	1.95	0.47
1:B:208:ASP:HB2	1:B:224:LYS:HE2	1.97	0.47
1:B:14:GLU:HA	1:B:17:LYS:HB2	1.97	0.47
1:B:141:GLU:O	1:B:145:ILE:HG13	2.15	0.47
1:B:31:GLN:HB3	1:B:31:GLN:HE21	1.60	0.47
1:A:160:LEU:HA	1:A:163:VAL:CG2	2.45	0.46
1:B:172:GLY:HA2	1:B:193:GLY:HA3	1.96	0.46
1:A:15:LYS:CD	1:A:64:ILE:HG13	2.46	0.46
1:B:132:ALA:O	1:B:135:VAL:HB	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:VAL:HG12	1:A:227:VAL:O	2.16	0.46
1:A:82:ASP:C	1:A:84:ILE:N	2.63	0.46
1:B:177:LEU:N	1:B:178:PRO:HD2	2.30	0.46
1:A:169:ALA:HA	1:A:190:TYR:OH	2.15	0.46
1:A:82:ASP:O	1:A:83:VAL:C	2.53	0.46
1:A:37:PRO:HA	1:A:93:LYS:HD3	1.96	0.46
1:B:153:PRO:HG3	1:B:246:PHE:CD2	2.50	0.46
1:A:124:GLY:O	1:A:125:ASN:O	2.33	0.46
1:A:221:LEU:O	1:A:222:ASP:HB3	2.15	0.46
1:A:31:GLN:O	1:A:34:LYS:NZ	2.48	0.46
1:A:52:PHE:CG	1:A:178:PRO:HG3	2.51	0.46
1:A:70:LEU:HA	1:A:70:LEU:HD12	1.64	0.46
1:A:20:ILE:HA	1:A:23:ILE:HG23	1.97	0.46
1:B:150:ARG:O	1:B:227:VAL:HG23	2.15	0.46
1:A:61:GLU:O	1:A:63:ALA:N	2.49	0.46
1:A:30:GLU:O	1:A:33:MET:HB2	2.16	0.46
1:A:154:ILE:HD12	1:A:159:TYR:HB2	1.98	0.46
1:B:22:GLN:O	1:B:26:GLN:HG2	2.15	0.45
1:A:274:LEU:O	1:A:278:VAL:HG23	2.16	0.45
1:A:227:VAL:HG21	1:A:246:PHE:CZ	2.50	0.45
1:A:70:LEU:HD23	1:A:170:LEU:HD13	1.99	0.45
1:A:35:ASP:O	1:A:93:LYS:HE3	2.17	0.45
1:B:102:ASP:O	1:B:105:ALA:HB3	2.16	0.45
1:A:29:PRO:HG2	1:B:148:LEU:CD1	2.46	0.45
1:A:233:VAL:O	1:A:234:THR:C	2.55	0.45
1:B:88:ARG:O	1:B:88:ARG:HG2	2.16	0.45
1:B:80:HIS:C	1:B:82:ASP:N	2.71	0.45
1:B:192:LEU:HD13	1:B:282:ASN:HD21	1.82	0.45
1:B:82:ASP:HB3	1:B:97:ASN:ND2	2.32	0.44
1:A:147:GLN:C	1:A:149:SER:N	2.71	0.44
1:B:201:GLN:O	1:B:204:ASP:HB2	2.17	0.44
1:A:105:ALA:O	1:A:108:ALA:HB3	2.18	0.44
1:B:83:VAL:HG13	1:B:97:ASN:HB2	2.00	0.44
1:A:230:PHE:N	1:A:231:PRO:HD2	2.32	0.44
1:B:28:PHE:HA	1:B:29:PRO:HD3	1.88	0.44
1:A:203:PHE:O	1:A:207:MET:HG2	2.18	0.44
1:A:29:PRO:O	1:A:32:ILE:HG22	2.18	0.44
1:A:102:ASP:O	1:A:106:VAL:HG23	2.18	0.44
1:A:202:MET:HB3	1:A:271:LEU:HD13	2.00	0.43
1:A:157:GLU:O	1:A:161:ARG:HG3	2.18	0.43
1:A:14:GLU:HA	1:A:17:LYS:HB2	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:THR:O	1:B:199:ILE:HG13	2.18	0.43
1:B:176:GLN:HG3	1:B:186:GLY:HA3	2.01	0.43
1:B:232:LEU:C	1:B:236:MET:HE2	2.37	0.43
1:A:116:ALA:O	1:A:119:THR:HB	2.19	0.43
1:B:27:PHE:CD2	1:B:115:SER:HB3	2.53	0.43
1:A:108:ALA:O	1:A:112:VAL:N	2.46	0.43
1:A:229:SER:HB3	1:A:232:LEU:HB3	1.99	0.43
1:A:230:PHE:CE2	1:A:267:CYS:HB3	2.54	0.43
1:A:276:LYS:C	1:A:280:ILE:HG12	2.38	0.43
1:B:78:LEU:HD23	1:B:78:LEU:HA	1.85	0.43
1:A:103:LYS:NZ	1:B:81:ASP:OD1	2.49	0.43
1:A:209:PHE:O	1:A:212:MET:HG2	2.19	0.43
1:A:229:SER:HB3	1:A:232:LEU:HB2	2.01	0.42
1:B:111:LEU:O	1:B:114:VAL:HB	2.18	0.42
1:A:112:VAL:O	1:A:113:LEU:C	2.57	0.42
1:A:19:ARG:O	1:A:23:ILE:CG2	2.66	0.42
1:A:21:GLU:HG3	1:A:33:MET:CE	2.48	0.42
1:A:38:LEU:O	1:A:39:TYR:O	2.38	0.42
1:B:170:LEU:HD12	1:B:170:LEU:HA	1.82	0.42
1:B:113:LEU:HG	1:B:117:PHE:HE1	1.84	0.42
1:B:73:VAL:HG11	1:B:136:ILE:HD12	2.01	0.42
1:B:288:PHE:CD1	1:B:288:PHE:C	2.92	0.42
1:B:213:GLU:HA	1:B:213:GLU:OE2	2.19	0.42
1:A:52:PHE:CD2	1:A:178:PRO:HG3	2.54	0.42
1:B:202:MET:HE2	1:B:202:MET:HB3	1.92	0.42
1:A:276:LYS:O	1:A:279:ILE:N	2.53	0.42
1:B:180:LEU:O	1:B:183:GLY:N	2.49	0.42
1:B:48:SER:O	1:B:52:PHE:HD2	2.03	0.42
1:A:241:GLU:HG3	1:A:242:ALA:N	2.34	0.42
1:A:82:ASP:O	1:A:84:ILE:N	2.53	0.42
1:A:21:GLU:O	1:A:25:SER:HB3	2.19	0.42
1:B:123:ILE:O	1:B:123:ILE:HG22	2.19	0.42
1:B:172:GLY:HA3	1:B:190:TYR:CD1	2.55	0.42
1:B:200:TYR:O	1:B:203:PHE:HB3	2.20	0.41
1:B:75:LEU:O	1:B:78:LEU:HB2	2.20	0.41
1:B:192:LEU:CD1	1:B:282:ASN:ND2	2.83	0.41
1:A:239:PHE:CD1	1:A:239:PHE:N	2.89	0.41
1:A:227:VAL:HG13	1:A:232:LEU:HD23	2.01	0.41
1:A:265:LYS:O	1:A:269:GLU:HG2	2.20	0.41
1:A:113:LEU:O	1:A:116:ALA:HB3	2.20	0.41
1:A:176:GLN:O	1:A:179:ALA:HB3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:ASN:O	1:A:194:VAL:HB	2.20	0.41
1:B:215:ILE:HG22	1:B:216:GLY:N	2.35	0.41
1:B:97:ASN:CA	1:B:105:ALA:HB2	2.49	0.41
1:A:114:VAL:O	1:A:114:VAL:HG12	2.21	0.41
1:A:149:SER:O	1:A:150:ARG:C	2.59	0.41
1:B:286:ARG:O	1:B:288:PHE:N	2.54	0.41
1:B:38:LEU:HA	1:B:38:LEU:HD23	1.87	0.41
1:B:229:SER:O	1:B:233:VAL:HG13	2.21	0.41
1:B:249:ARG:NE	1:B:251:TRP:HH2	2.19	0.40
1:A:69:ALA:HA	1:A:72:LEU:HD12	2.02	0.40
1:A:176:GLN:NE2	1:A:190:TYR:HB2	2.35	0.40
1:B:88:ARG:HG3	1:B:98:PHE:CD2	2.56	0.40
1:A:55:ARG:NE	1:A:179:ALA:HA	2.36	0.40
1:A:146:GLU:OE2	1:A:225:ASN:HB3	2.22	0.40
1:A:175:LEU:O	1:A:178:PRO:HD2	2.21	0.40
1:A:283:SER:HA	1:A:286:ARG:NH2	2.35	0.40
1:B:275:VAL:O	1:B:278:VAL:HB	2.21	0.40
1:B:57:VAL:CG1	1:B:58:GLU:N	2.85	0.40
1:A:144:LEU:O	1:A:148:LEU:HG	2.22	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:476:HOH:O	2:A:476:HOH:O[7_556]	0.98	1.22
2:A:409:HOH:O	2:A:410:HOH:O[2_555]	1.99	0.21

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	278/299 (93%)	228 (82%)	41 (15%)	9 (3%)	5 31

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	278/299 (93%)	231 (83%)	34 (12%)	13 (5%)	3	20
All	All	556/598 (93%)	459 (83%)	75 (14%)	22 (4%)	4	24

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	39	TYR
1	A	61	GLU
1	B	38	LEU
1	B	187	GLU
1	B	238	LYS
1	A	125	ASN
1	B	61	GLU
1	B	221	LEU
1	A	59	ILE
1	A	150	ARG
1	A	280	ILE
1	B	121	GLU
1	B	81	ASP
1	B	95	THR
1	B	150	ARG
1	B	283	SER
1	A	38	LEU
1	A	221	LEU
1	B	91	ARG
1	B	210	ALA
1	B	287	ASP
1	A	112	VAL

5.3.2 Protein sidechains ⓘ

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	237/255 (93%)	220 (93%)	17 (7%)	18	54
1	B	237/255 (93%)	225 (95%)	12 (5%)	29	68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	474/510 (93%)	445 (94%)	29 (6%)	23	62

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	TYR
1	A	23	ILE
1	A	26	GLN
1	A	28	PHE
1	A	30	GLU
1	A	39	TYR
1	A	50	LEU
1	A	66	SER
1	A	75	LEU
1	A	81	ASP
1	A	85	ASP
1	A	157	GLU
1	A	204	ASP
1	A	218	ASP
1	A	237	GLU
1	A	241	GLU
1	A	259	ARG
1	B	30	GLU
1	B	31	GLN
1	B	34	LYS
1	B	62	ASP
1	B	65	SER
1	B	81	ASP
1	B	83	VAL
1	B	110	ASP
1	B	136	ILE
1	B	188	ASP
1	B	204	ASP
1	B	229	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	54	ASN
1	A	80	HIS

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Mol	Chain	Res	Type
1	A	147	GLN
1	A	191	ASN
1	A	225	ASN
1	A	282	ASN
1	B	31	GLN
1	B	191	ASN
1	B	248	ASN
1	B	282	ASN

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	280/299 (93%)	0.41	22 (7%) 15 12	41, 67, 94, 106	0
1	B	280/299 (93%)	0.31	15 (5%) 29 24	27, 62, 88, 106	0
All	All	560/598 (93%)	0.36	37 (6%) 22 17	27, 64, 91, 106	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	153	PRO	5.2
1	B	242	ALA	4.9
1	B	148	LEU	4.6
1	B	246	PHE	4.4
1	A	123	ILE	4.1
1	A	64	ILE	4.0
1	A	181	LEU	4.0
1	A	122	GLU	4.0
1	B	124	GLY	3.4
1	A	246	PHE	3.3
1	A	285	LEU	3.1
1	B	151	TYR	3.0
1	A	65	SER	3.0
1	A	59	ILE	2.9
1	A	242	ALA	2.8
1	A	61	GLU	2.8
1	A	183	GLY	2.8
1	B	95	THR	2.6
1	A	124	GLY	2.5
1	A	239	PHE	2.4
1	B	106	VAL	2.4
1	A	153	PRO	2.3
1	B	181	LEU	2.3
1	A	67	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	285	LEU	2.2
1	B	241	GLU	2.2
1	B	248	ASN	2.2
1	A	286	ARG	2.2
1	B	187	GLU	2.2
1	A	241	GLU	2.2
1	A	248	ASN	2.2
1	B	244	GLN	2.1
1	A	180	LEU	2.1
1	A	129	ARG	2.1
1	A	57	VAL	2.0
1	A	9	ASN	2.0
1	B	287	ASP	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.