



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 12:15 am BST

PDB ID : 3AG5  
Title : Crystal Structure of Pantothenate Synthetase from *Staphylococcus aureus*  
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Deposited on : 2010-03-19  
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](mailto: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.

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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

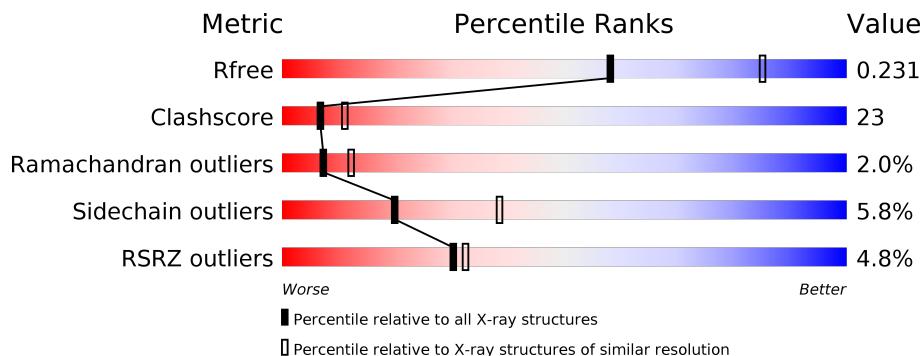
# 1 Overall quality at a glance

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  $\geq 3$ , 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	283	 3% 66% 28%
1	B	283	 7% 55% 37% 6%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4404 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 Pantothenate synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	278	Total 2178	C 1387	N 368	O 416	S 7	0	0	0
1	B	279	Total 2193	C 1398	N 369	O 419	S 7	0	0	0

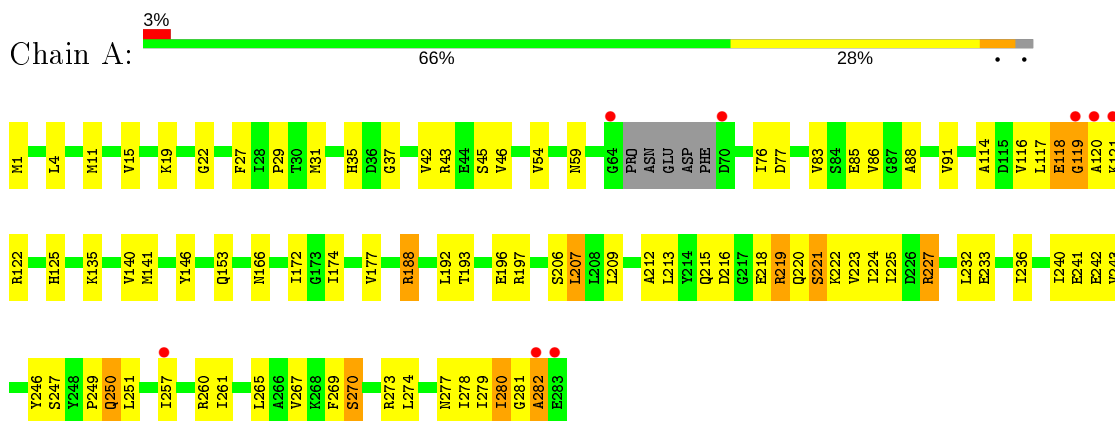
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	18	Total 18	O 18	0	0
2	B	15	Total 15	O 15	0	0

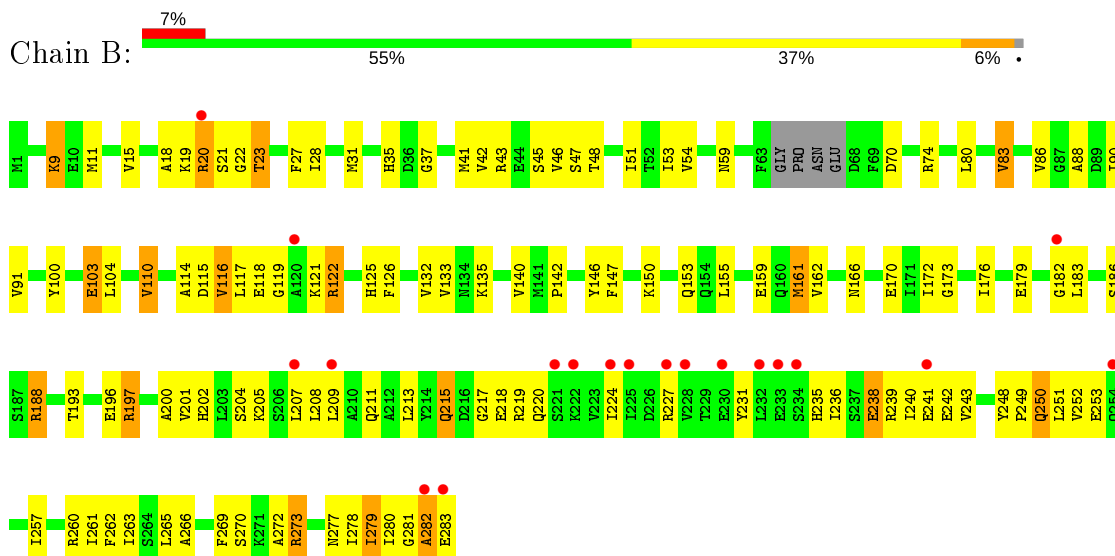
### 3 Residue-property plots [i](#)

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: Pantothenate synthetase



- Molecule 1: Pantothenate synthetase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	131.66 Å 131.66 Å 89.24 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	41.55 – 2.50 41.55 – 2.49	Depositor EDS
% Data completeness (in resolution range)	99.4 (41.55-2.50) 99.1 (41.55-2.49)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.38 (at 2.48 Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.221 , 0.241 0.221 , 0.231	Depositor DCC
$R_{free}$ test set	1535 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	58.2	Xtrriage
Anisotropy	0.130	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 39.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.039 for h,-h-k,-l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4404	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	71.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.08% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.39	0/2213	0.64	0/2993
1	B	0.38	0/2229	0.62	0/3015
All	All	0.38	0/4442	0.63	0/6008

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 [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	2178	0	2207	83	0
1	B	2193	0	2217	130	0
2	A	18	0	0	0	0
2	B	15	0	0	0	0
All	All	4404	0	4424	205	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 23.

All (205) 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:213:LEU:HD11	1:B:227:ARG:HG2	1.34	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:ARG:HG3	1:A:257:ILE:HB	1.43	0.98
1:B:110:VAL:HG22	1:B:114:ALA:HB2	1.50	0.94
1:A:114:ALA:O	1:A:119:GLY:HA3	1.68	0.93
1:B:188:ARG:HB3	1:B:188:ARG:HH11	1.36	0.91
1:B:273:ARG:HH11	1:B:273:ARG:CG	1.85	0.90
1:A:31:MET:CE	1:A:59:ASN:H	1.86	0.87
1:B:238:GLU:HG2	1:B:269:PHE:HA	1.58	0.86
1:B:273:ARG:HG3	1:B:273:ARG:HH11	1.40	0.86
1:B:273:ARG:NH1	1:B:273:ARG:HG3	1.91	0.85
1:B:11:MET:HE1	1:B:53:ILE:HD13	1.57	0.85
1:B:83:VAL:HG13	1:B:88:ALA:HB3	1.61	0.83
1:B:11:MET:CE	1:B:53:ILE:HD13	2.12	0.80
1:B:31:MET:CE	1:B:59:ASN:H	1.95	0.80
1:B:35:HIS:HD2	1:B:37:GLY:H	1.27	0.79
1:A:257:ILE:HD12	1:A:261:ILE:HG21	1.64	0.79
1:B:211:GLN:HE21	1:B:280:ILE:HG21	1.49	0.77
1:A:83:VAL:HG22	1:A:88:ALA:HB2	1.67	0.77
1:B:239:ARG:HB3	1:B:239:ARG:NH1	2.02	0.75
1:B:219:ARG:HH21	1:B:219:ARG:HG3	1.50	0.75
1:B:31:MET:HE2	1:B:59:ASN:H	1.54	0.72
1:A:116:VAL:HG21	1:A:251:LEU:O	1.89	0.72
1:A:166:ASN:HD21	1:B:135:LYS:NZ	1.88	0.71
1:A:135:LYS:HZ1	1:B:166:ASN:HD21	1.39	0.70
1:A:35:HIS:HD2	1:A:37:GLY:H	1.38	0.70
1:B:188:ARG:HH11	1:B:188:ARG:CB	2.05	0.70
1:A:250:GLN:NE2	1:A:250:GLN:H	1.89	0.70
1:B:9:LYS:C	1:B:9:LYS:HD3	2.12	0.70
1:A:196:GLU:HB2	1:A:269:PHE:CD2	2.27	0.69
1:B:202:HIS:CE1	1:B:235:HIS:HB3	2.28	0.68
1:A:220:GLN:O	1:A:223:VAL:HG22	1.94	0.68
1:B:220:GLN:O	1:B:224:ILE:HG12	1.93	0.68
1:A:218:GLU:CD	1:A:223:VAL:HG21	2.14	0.68
1:B:239:ARG:HB3	1:B:239:ARG:HH11	1.57	0.66
1:A:31:MET:HE2	1:A:59:ASN:H	1.60	0.66
1:A:135:LYS:NZ	1:B:166:ASN:HD21	1.93	0.66
1:A:218:GLU:OE2	1:A:223:VAL:HG21	1.96	0.65
1:B:262:PHE:HD2	1:B:279:ILE:HG23	1.61	0.65
1:B:19:LYS:C	1:B:21:SER:H	1.98	0.65
1:A:15:VAL:CG1	1:A:141:MET:HB2	2.28	0.64
1:B:21:SER:O	1:B:23:THR:N	2.25	0.64
1:A:15:VAL:HG11	1:A:141:MET:HB2	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:ARG:HE	1:B:125:HIS:CG	2.16	0.63
1:A:232:LEU:O	1:A:236:ILE:HG12	1.98	0.63
1:A:207:LEU:HD13	1:A:265:LEU:HD11	1.80	0.63
1:A:219:ARG:CG	1:A:257:ILE:HB	2.24	0.63
1:B:59:ASN:CG	1:B:132:VAL:HG11	2.18	0.62
1:B:236:ILE:HD11	1:B:240:ILE:HD11	1.82	0.62
1:B:42:VAL:O	1:B:46:VAL:HG23	2.00	0.62
1:A:233:GLU:HG3	1:A:240:ILE:HD12	1.83	0.61
1:B:260:ARG:HD2	1:B:283:GLU:OXT	2.00	0.61
1:A:54:VAL:HG21	1:A:83:VAL:HG21	1.80	0.61
1:A:250:GLN:CD	1:A:250:GLN:H	2.04	0.61
1:A:221:SER:O	1:A:225:ILE:HG13	2.01	0.60
1:B:155:LEU:O	1:B:159:GLU:HG3	2.01	0.60
1:B:188:ARG:HB3	1:B:188:ARG:NH1	2.13	0.60
1:B:202:HIS:ND1	1:B:235:HIS:HB3	2.15	0.60
1:A:19:LYS:HE2	1:A:141:MET:O	2.01	0.60
1:A:213:LEU:HD11	1:A:227:ARG:HG2	1.82	0.60
1:A:193:THR:OG1	1:A:196:GLU:HG2	2.02	0.60
1:A:118:GLU:HG3	1:A:119:GLY:N	2.17	0.59
1:B:150:LYS:HB2	1:B:176:ILE:HG22	1.84	0.59
1:A:35:HIS:CD2	1:A:37:GLY:H	2.20	0.59
1:A:45:SER:HB2	1:A:146:TYR:CE2	2.38	0.59
1:B:179:GLU:HG2	1:B:197:ARG:HH12	1.69	0.58
1:A:11:MET:HE2	1:A:15:VAL:HG23	1.84	0.58
1:B:83:VAL:HG13	1:B:88:ALA:CB	2.33	0.58
1:A:222:LYS:O	1:A:222:LYS:HD3	2.04	0.58
1:A:86:VAL:O	1:A:86:VAL:HG12	2.04	0.57
1:A:166:ASN:HD21	1:B:135:LYS:HZ3	1.52	0.57
1:B:179:GLU:HG2	1:B:197:ARG:NH1	2.19	0.57
1:A:31:MET:HE1	1:A:59:ASN:H	1.69	0.57
1:A:196:GLU:HB2	1:A:269:PHE:HD2	1.69	0.57
1:B:86:VAL:HG12	1:B:86:VAL:O	2.04	0.57
1:A:22:GLY:HA2	1:B:20:ARG:HD2	1.87	0.57
1:B:219:ARG:HG3	1:B:219:ARG:NH2	2.19	0.57
1:A:257:ILE:HD12	1:A:261:ILE:CG2	2.33	0.56
1:A:135:LYS:NZ	1:B:166:ASN:ND2	2.53	0.56
1:B:236:ILE:HD11	1:B:240:ILE:CD1	2.36	0.56
1:B:239:ARG:HH11	1:B:239:ARG:CB	2.19	0.56
1:B:116:VAL:HG13	1:B:117:LEU:HG	1.88	0.55
1:B:207:LEU:HD13	1:B:265:LEU:HD11	1.89	0.55
1:B:54:VAL:O	1:B:91:VAL:HA	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:GLN:NE2	1:B:280:ILE:HG21	2.20	0.55
1:A:207:LEU:CD1	1:A:265:LEU:HD11	2.37	0.54
1:B:103:GLU:OE1	1:B:103:GLU:HA	2.07	0.54
1:B:241:GLU:HB3	1:B:266:ALA:O	2.07	0.53
1:A:192:LEU:HD11	1:A:274:LEU:HD11	1.91	0.53
1:A:85:GLU:HG2	1:A:85:GLU:O	2.07	0.53
1:A:227:ARG:HG3	1:A:227:ARG:HH11	1.74	0.52
1:A:54:VAL:O	1:A:91:VAL:HA	2.09	0.52
1:B:115:ASP:HA	1:B:119:GLY:HA3	1.90	0.52
1:B:207:LEU:HD13	1:B:265:LEU:CD1	2.40	0.52
1:A:215:GLN:HE21	1:A:215:GLN:HA	1.75	0.52
1:B:122:ARG:HH21	1:B:122:ARG:HG3	1.74	0.52
1:B:218:GLU:HG3	1:B:224:ILE:HG23	1.92	0.52
1:A:11:MET:CE	1:A:15:VAL:HG23	2.40	0.52
1:A:193:THR:HG23	1:A:196:GLU:OE1	2.10	0.51
1:A:83:VAL:HG22	1:A:88:ALA:CB	2.39	0.51
1:B:238:GLU:CG	1:B:269:PHE:HA	2.36	0.51
1:A:212:ALA:O	1:A:215:GLN:HB3	2.09	0.51
1:B:269:PHE:HB2	1:B:272:ALA:O	2.11	0.51
1:A:116:VAL:HG22	1:A:117:LEU:HG	1.93	0.51
1:A:118:GLU:O	1:A:120:ALA:N	2.43	0.51
1:B:35:HIS:CD2	1:B:37:GLY:H	2.18	0.50
1:B:41:MET:HE1	1:B:147:PHE:HA	1.93	0.50
1:B:43:ARG:HG3	1:B:86:VAL:CG1	2.42	0.50
1:A:118:GLU:HB3	1:A:153:GLN:OE1	2.12	0.50
1:A:166:ASN:HD21	1:B:135:LYS:HZ2	1.59	0.50
1:B:227:ARG:HD2	1:B:227:ARG:O	2.12	0.50
1:B:257:ILE:HG21	1:B:261:ILE:HG21	1.94	0.50
1:A:116:VAL:HG22	1:A:117:LEU:CD2	2.41	0.50
1:B:70:ASP:HA	1:B:74:ARG:HH21	1.77	0.49
1:A:15:VAL:HG11	1:A:140:VAL:O	2.12	0.49
1:A:11:MET:HE2	1:A:15:VAL:CG2	2.42	0.49
1:B:118:GLU:HG3	1:B:118:GLU:O	2.13	0.49
1:B:205:LYS:C	1:B:207:LEU:H	2.15	0.49
1:B:218:GLU:HG3	1:B:224:ILE:CG2	2.42	0.49
1:B:182:GLY:O	1:B:204:SER:HB3	2.12	0.49
1:B:118:GLU:HB2	1:B:121:LYS:HD3	1.94	0.49
1:B:186:SER:OG	1:B:188:ARG:HG3	2.13	0.48
1:B:19:LYS:C	1:B:21:SER:N	2.66	0.48
1:B:147:PHE:O	1:B:173:GLY:HA2	2.13	0.48
1:A:261:ILE:O	1:A:279:ILE:HA	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:201:VAL:O	1:B:201:VAL:HG12	2.14	0.48
1:B:150:LYS:HB2	1:B:176:ILE:CG2	2.44	0.47
1:B:183:LEU:HA	1:B:204:SER:OG	2.13	0.47
1:A:241:GLU:O	1:A:242:GLU:HB2	2.14	0.47
1:B:28:ILE:HD12	1:B:28:ILE:N	2.29	0.47
1:B:80:LEU:O	1:B:80:LEU:HD23	2.14	0.47
1:A:166:ASN:ND2	1:B:135:LYS:NZ	2.59	0.47
1:A:118:GLU:HG3	1:A:119:GLY:H	1.78	0.47
1:A:236:ILE:HD12	1:A:267:VAL:HG11	1.97	0.47
1:B:242:GLU:HG2	1:B:243:VAL:N	2.29	0.46
1:B:248:TYR:CE1	1:B:249:PRO:HG3	2.50	0.46
1:B:215:GLN:C	1:B:215:GLN:HE21	2.18	0.46
1:A:122:ARG:HG2	1:A:125:HIS:HB2	1.97	0.46
1:A:277:ASN:O	1:A:278:ILE:HG23	2.15	0.46
1:B:186:SER:OG	1:B:188:ARG:CG	2.64	0.46
1:B:9:LYS:O	1:B:9:LYS:HD3	2.14	0.46
1:B:132:VAL:HG23	1:B:133:VAL:N	2.31	0.46
1:A:223:VAL:HG23	1:A:224:ILE:N	2.29	0.46
1:B:277:ASN:O	1:B:278:ILE:HG23	2.15	0.45
1:A:206:SER:O	1:A:209:LEU:HB3	2.17	0.45
1:A:273:ARG:HG3	1:A:273:ARG:O	2.16	0.45
1:B:19:LYS:O	1:B:21:SER:N	2.49	0.45
1:B:262:PHE:CD2	1:B:279:ILE:HG23	2.48	0.45
1:A:233:GLU:HG3	1:A:240:ILE:CD1	2.45	0.45
1:B:205:LYS:HA	1:B:208:LEU:HG	1.99	0.45
1:A:269:PHE:O	1:A:270:SER:C	2.55	0.45
1:B:273:ARG:O	1:B:273:ARG:CG	2.64	0.45
1:A:281:GLY:O	1:A:282:ALA:HB2	2.16	0.45
1:B:202:HIS:HD2	1:B:205:LYS:NZ	2.14	0.45
1:B:250:GLN:HB2	1:B:252:VAL:HG22	1.99	0.44
1:B:15:VAL:HG11	1:B:140:VAL:O	2.17	0.44
1:B:45:SER:HB2	1:B:146:TYR:CE2	2.53	0.44
1:A:236:ILE:HG13	1:A:236:ILE:O	2.16	0.44
1:B:261:ILE:O	1:B:279:ILE:HA	2.18	0.44
1:B:161:MET:HE3	1:B:162:VAL:N	2.33	0.44
1:B:207:LEU:HG	1:B:278:ILE:HG12	1.99	0.44
1:B:53:ILE:HG12	1:B:90:ILE:HB	2.00	0.44
1:B:213:LEU:HB3	1:B:218:GLU:HB3	2.00	0.43
1:A:42:VAL:O	1:A:46:VAL:HG23	2.17	0.43
1:B:273:ARG:HG2	1:B:273:ARG:O	2.18	0.43
1:B:281:GLY:O	1:B:282:ALA:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:ARG:HG2	1:B:282:ALA:H	1.84	0.43
1:B:146:TYR:CD1	1:B:172:ILE:HB	2.53	0.43
1:A:76:ILE:HG23	1:A:77:ASP:N	2.32	0.43
1:B:47:SER:OG	1:B:48:THR:HG23	2.18	0.43
1:B:15:VAL:O	1:B:19:LYS:HG3	2.19	0.43
1:B:21:SER:O	1:B:23:THR:HG22	2.18	0.43
1:B:217:GLY:O	1:B:218:GLU:C	2.56	0.43
1:A:27:PHE:CE2	1:A:29:PRO:HG3	2.53	0.43
1:A:250:GLN:NE2	1:A:250:GLN:N	2.62	0.42
1:B:146:TYR:HA	1:B:172:ILE:O	2.18	0.42
1:B:263:ILE:O	1:B:277:ASN:HB2	2.18	0.42
1:B:193:THR:OG1	1:B:196:GLU:HG3	2.18	0.42
1:B:27:PHE:HB2	1:B:142:PRO:HB3	2.01	0.42
1:B:205:LYS:C	1:B:207:LEU:N	2.72	0.42
1:A:247:SER:O	1:A:251:LEU:HA	2.19	0.42
1:B:116:VAL:HG21	1:B:251:LEU:HB2	2.02	0.42
1:B:200:ALA:C	1:B:202:HIS:H	2.23	0.42
1:B:110:VAL:CG2	1:B:114:ALA:HB2	2.36	0.42
1:A:117:LEU:HD21	1:A:246:TYR:CE2	2.55	0.41
1:A:249:PRO:O	1:A:251:LEU:HG	2.19	0.41
1:B:209:LEU:HG	1:B:227:ARG:HH21	1.85	0.41
1:B:205:LYS:HZ2	1:B:205:LYS:HB2	1.86	0.41
1:B:238:GLU:OE1	1:B:238:GLU:N	2.54	0.41
1:B:18:ALA:O	1:B:23:THR:CG2	2.68	0.41
1:A:188:ARG:HH11	1:A:188:ARG:HG2	1.86	0.41
1:B:161:MET:HE2	1:B:161:MET:HB3	1.97	0.41
1:B:116:VAL:HG21	1:B:251:LEU:CB	2.50	0.41
1:B:261:ILE:O	1:B:279:ILE:HG22	2.20	0.41
1:A:280:ILE:N	1:A:280:ILE:HD13	2.36	0.41
1:B:126:PHE:CZ	1:B:153:GLN:HG2	2.56	0.41
1:B:253:GLU:O	1:B:253:GLU:HG3	2.20	0.41
1:A:172:ILE:HG22	1:A:174:ILE:CD1	2.50	0.41
1:B:100:TYR:OH	1:B:132:VAL:HG12	2.20	0.41
1:B:205:LYS:NZ	1:B:205:LYS:CB	2.84	0.41
1:B:51:ILE:O	1:B:51:ILE:HG13	2.20	0.41
1:A:43:ARG:HG3	1:A:86:VAL:CG1	2.51	0.40
1:B:231:TYR:C	1:B:231:TYR:CD1	2.94	0.40
1:B:100:TYR:OH	1:B:132:VAL:CG1	2.69	0.40
1:A:119:GLY:C	1:A:121:LYS:H	2.25	0.40
1:B:41:MET:CE	1:B:147:PHE:C	2.90	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	274/283 (97%)	249 (91%)	19 (7%)	6 (2%)	6	10
1	B	275/283 (97%)	244 (89%)	26 (10%)	5 (2%)	8	14
All	All	549/566 (97%)	493 (90%)	45 (8%)	11 (2%)	7	12

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	118	GLU
1	A	119	GLY
1	A	282	ALA
1	B	22	GLY
1	B	270	SER
1	B	282	ALA
1	A	270	SER
1	B	20	ARG
1	A	221	SER
1	B	122	ARG
1	A	243	VAL

### 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	239/244 (98%)	227 (95%)	12 (5%)	24	46
1	B	241/244 (99%)	225 (93%)	16 (7%)	16	32

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	480/488 (98%)	452 (94%)	28 (6%)	20	38

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	4	LEU
1	A	177	VAL
1	A	188	ARG
1	A	197	ARG
1	A	207	LEU
1	A	216	ASP
1	A	219	ARG
1	A	227	ARG
1	A	250	GLN
1	A	260	ARG
1	A	280	ILE
1	B	9	LYS
1	B	23	THR
1	B	83	VAL
1	B	103	GLU
1	B	104	LEU
1	B	110	VAL
1	B	116	VAL
1	B	161	MET
1	B	170	GLU
1	B	188	ARG
1	B	197	ARG
1	B	215	GLN
1	B	238	GLU
1	B	250	GLN
1	B	273	ARG
1	B	279	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	HIS
1	A	49	ASN
1	A	125	HIS
1	A	166	ASN

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Mol	Chain	Res	Type
1	A	211	GLN
1	A	215	GLN
1	A	250	GLN
1	A	254	GLN
1	B	35	HIS
1	B	49	ASN
1	B	125	HIS
1	B	166	ASN
1	B	198	GLN
1	B	202	HIS
1	B	211	GLN
1	B	215	GLN
1	B	235	HIS
1	B	254	GLN

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	278/283 (98%)	-0.04	8 (2%) 51 55	39, 60, 98, 120	0
1	B	279/283 (98%)	0.17	19 (6%) 17 17	41, 71, 125, 129	0
All	All	557/566 (98%)	0.07	27 (4%) 30 32	39, 65, 118, 129	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	120	ALA	4.4
1	B	232	LEU	4.2
1	A	282	ALA	4.2
1	B	283	GLU	3.9
1	A	283	GLU	3.5
1	B	254	GLN	3.1
1	A	121	LYS	2.9
1	B	282	ALA	2.8
1	B	233	GLU	2.7
1	B	224	ILE	2.7
1	B	230	GLU	2.6
1	B	222	LYS	2.5
1	A	119	GLY	2.5
1	A	70	ASP	2.5
1	B	227	ARG	2.4
1	B	221	SER	2.4
1	A	257	ILE	2.4
1	B	209	LEU	2.4
1	B	120	ALA	2.3
1	B	228	VAL	2.3
1	B	225	ILE	2.3
1	B	20	ARG	2.2
1	B	207	LEU	2.2
1	B	234	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	64	GLY	2.1
1	B	182	GLY	2.1
1	B	241	GLU	2.1

## 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.