



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:40 PM GMT

PDB ID : 3RMG  
Title : Crystal structure of geranylgeranyl pyrophosphate synthase from bacteroides thetaiotaomicron  
Authors : Patskovsky, Y.; Toro, R.; Sauder, J.M.; Poulter, C.D.; Gerlt, J.A.; Burley, S.K.; Almo, S.C.; Enzyme Function Initiative (EFI); New York SGX Research Center for Structural Genomics (NYSGXRC)  
Deposited on : 2011-04-20  
Resolution : 2.30 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](mailto: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.

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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  
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

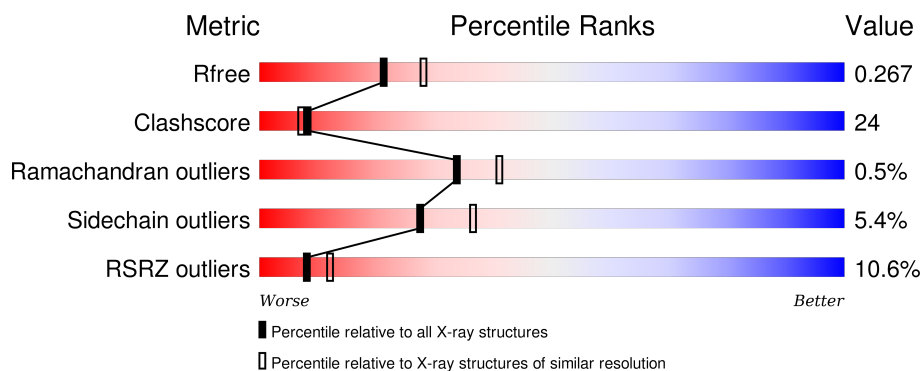
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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. 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	334	
1	B	334	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4627 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 Octaprenyl-diphosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	299	Total	C	N	O	S	0	1	0
			2301	1472	369	453	7			
1	B	297	Total	C	N	O	S	0	2	0
			2257	1444	367	439	7			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	MET	-	EXPRESSION TAG	UNP Q8A2P4
A	-1	SER	-	EXPRESSION TAG	UNP Q8A2P4
A	0	LEU	-	EXPRESSION TAG	UNP Q8A2P4
A	325	GLU	-	EXPRESSION TAG	UNP Q8A2P4
A	326	GLY	-	EXPRESSION TAG	UNP Q8A2P4
A	327	HIS	-	EXPRESSION TAG	UNP Q8A2P4
A	328	HIS	-	EXPRESSION TAG	UNP Q8A2P4
A	329	HIS	-	EXPRESSION TAG	UNP Q8A2P4
A	330	HIS	-	EXPRESSION TAG	UNP Q8A2P4
A	331	HIS	-	EXPRESSION TAG	UNP Q8A2P4
A	332	HIS	-	EXPRESSION TAG	UNP Q8A2P4
B	-2	MET	-	EXPRESSION TAG	UNP Q8A2P4
B	-1	SER	-	EXPRESSION TAG	UNP Q8A2P4
B	0	LEU	-	EXPRESSION TAG	UNP Q8A2P4
B	325	GLU	-	EXPRESSION TAG	UNP Q8A2P4
B	326	GLY	-	EXPRESSION TAG	UNP Q8A2P4
B	327	HIS	-	EXPRESSION TAG	UNP Q8A2P4
B	328	HIS	-	EXPRESSION TAG	UNP Q8A2P4
B	329	HIS	-	EXPRESSION TAG	UNP Q8A2P4
B	330	HIS	-	EXPRESSION TAG	UNP Q8A2P4
B	331	HIS	-	EXPRESSION TAG	UNP Q8A2P4
B	332	HIS	-	EXPRESSION TAG	UNP Q8A2P4

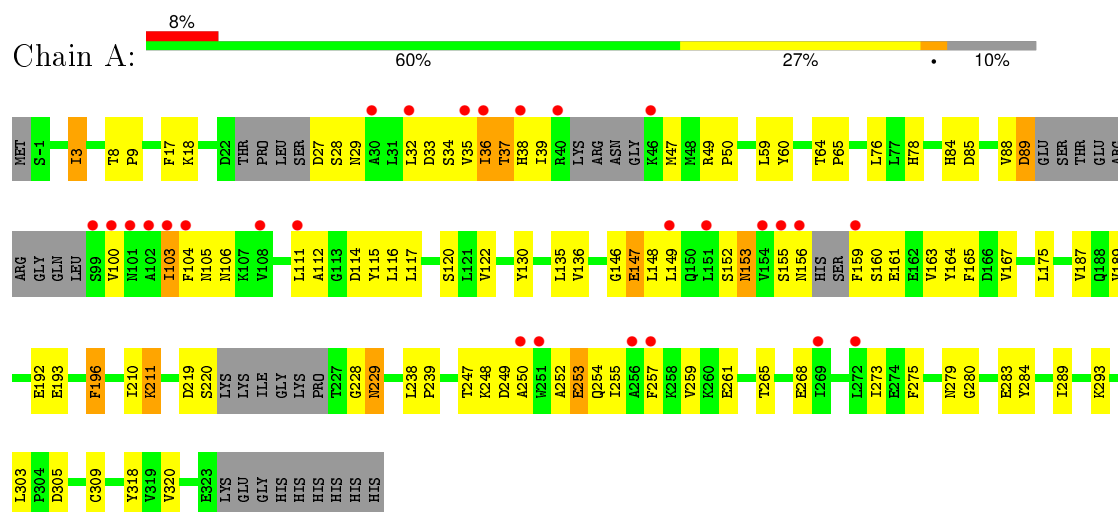
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	43	Total 43	O 43	0	0
2	B	26	Total 26	O 26	0	0

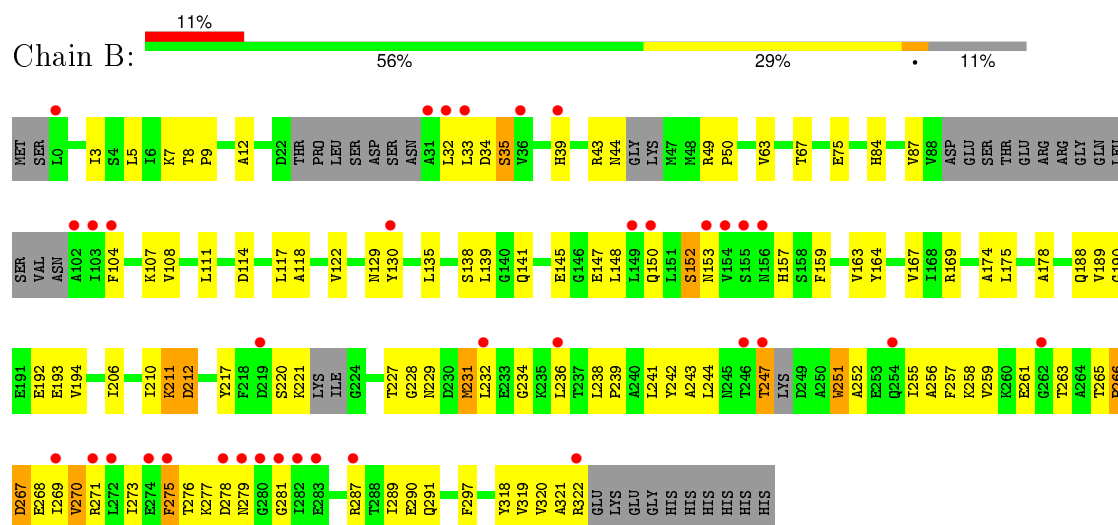
### 3 Residue-property plots

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



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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	44.62Å 111.41Å 83.38Å 90.00° 92.56° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30 44.57 – 2.19	Depositor EDS
% Data completeness (in resolution range)	99.7 (50.00-2.30) 99.2 (44.57-2.19)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.48 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.214 , 0.258 0.217 , 0.267	Depositor DCC
$R_{free}$ test set	1122 reflections (3.21%)	DCC
Wilson B-factor (Å <sup>2</sup> )	46.5	Xtriage
Anisotropy	0.318	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 64.4	EDS
Estimated twinning fraction	0.037 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 41295 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4627	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.31% 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, 0.375 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	0.48	0/2335	0.59	0/3168
1	B	0.47	0/2295	0.53	0/3119
All	All	0.48	0/4630	0.56	0/6287

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	2301	0	2283	112	0
1	B	2257	0	2216	123	0
2	A	43	0	0	0	0
2	B	26	0	0	3	0
All	All	4627	0	4499	221	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 24.

All (221) 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:36:ILE:HD12	1:A:36:ILE:C	1.55	1.21

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:LEU:HD22	1:B:145:GLU:HG3	1.25	1.15
1:A:33:ASP:O	1:A:37:THR:HB	1.46	1.15
1:A:88:VAL:HG13	1:A:89:ASP:OD1	1.49	1.12
1:B:49:ARG:HD3	1:B:175:LEU:HD12	1.25	1.07
1:A:103:ILE:O	1:A:103:ILE:HD12	1.56	1.04
1:A:275:PHE:CE1	1:A:279:ASN:ND2	2.30	0.99
1:B:269:ILE:O	1:B:273:ILE:HG13	1.62	0.99
1:A:36:ILE:HD12	1:A:36:ILE:O	1.61	0.98
1:A:249:ASP:O	1:A:253:GLU:HG3	1.64	0.96
1:A:255:ILE:HD12	1:A:268:GLU:OE2	1.65	0.96
1:A:36:ILE:C	1:A:36:ILE:CD1	2.29	0.93
1:A:32:LEU:O	1:A:36:ILE:HG23	1.67	0.93
1:B:251:TRP:CD1	1:B:255:ILE:HD11	2.04	0.91
1:A:293:LYS:HD2	1:A:320:VAL:HG21	1.52	0.90
1:B:49:ARG:HD3	1:B:175:LEU:CD1	1.99	0.90
1:A:117:LEU:HD23	1:B:117:LEU:HD23	1.54	0.89
1:B:243:ALA:HB1	1:B:275:PHE:CZ	2.08	0.88
1:A:36:ILE:HD12	1:A:37:THR:N	1.89	0.87
1:A:85:ASP:O	1:A:88:VAL:HG12	1.75	0.86
1:A:29:ASN:HB3	1:A:32:LEU:HB3	1.57	0.86
1:A:103:ILE:C	1:A:103:ILE:HD12	1.96	0.85
1:B:267:ASP:O	1:B:270:VAL:HB	1.76	0.85
1:A:29:ASN:HB3	1:A:32:LEU:CB	2.09	0.83
1:B:227:THR:HG22	1:B:273:ILE:CD1	2.09	0.83
1:B:104:PHE:O	1:B:108:VAL:HG23	1.78	0.82
1:A:146:GLY:O	1:A:149:LEU:HB3	1.82	0.79
1:A:34:SER:O	1:A:38:HIS:HB2	1.85	0.77
1:A:29:ASN:CB	1:A:32:LEU:HB3	2.14	0.77
1:A:103:ILE:HG13	1:A:104:PHE:N	1.98	0.76
1:B:251:TRP:CD1	1:B:252:ALA:N	2.55	0.75
1:B:227:THR:CG2	1:B:273:ILE:HD13	2.17	0.75
1:B:232:LEU:O	1:B:232:LEU:HD12	1.87	0.74
1:B:231:MET:O	1:B:259:VAL:HG13	1.86	0.74
1:A:255:ILE:CD1	1:A:268:GLU:OE2	2.37	0.72
1:B:243:ALA:CB	1:B:275:PHE:CZ	2.73	0.71
1:B:270:VAL:HG12	1:B:271:ARG:N	2.06	0.70
1:A:76:LEU:CB	1:A:120[A]:SER:OG	2.39	0.69
1:B:251:TRP:HD1	1:B:255:ILE:HD11	1.57	0.69
1:B:227:THR:HG22	1:B:273:ILE:HD13	1.73	0.69
1:B:148:LEU:CD1	1:B:148:LEU:N	2.56	0.68
1:B:251:TRP:C	1:B:251:TRP:CD1	2.65	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:273:ILE:HG22	1:B:277:LYS:HE3	1.75	0.67
1:A:117:LEU:HD23	1:B:117:LEU:CD2	2.24	0.67
1:B:273:ILE:O	1:B:277:LYS:HG3	1.94	0.67
1:B:211:LYS:HB3	1:B:322:ARG:HH12	1.59	0.67
1:A:280:GLY:HA2	1:A:283:GLU:OE1	1.96	0.66
1:A:106:ASN:HD22	1:A:106:ASN:N	1.91	0.66
1:B:163:VAL:O	1:B:167:VAL:HG23	1.94	0.66
1:A:275:PHE:CD1	1:A:279:ASN:ND2	2.63	0.65
1:A:36:ILE:CD1	1:A:37:THR:N	2.53	0.65
1:B:321:ALA:O	1:B:322:ARG:C	2.36	0.64
1:B:148:LEU:HD12	1:B:148:LEU:N	2.12	0.64
1:B:236:LEU:O	1:B:241:LEU:HG	1.98	0.64
1:B:164:TYR:CE2	1:B:238:LEU:HB2	2.33	0.64
1:B:239:PRO:HB3	1:B:281:GLY:O	1.99	0.63
1:A:29:ASN:CB	1:A:32:LEU:CB	2.74	0.63
1:A:229:ASN:HD22	1:A:229:ASN:C	1.99	0.62
1:A:76:LEU:HB3	1:A:120[A]:SER:OG	2.00	0.62
1:A:228:GLY:HA2	1:A:273:ILE:HD11	1.81	0.62
1:B:227:THR:CG2	1:B:273:ILE:CD1	2.75	0.62
1:B:43:ARG:O	1:B:44:ASN:C	2.38	0.61
1:A:147:GLU:HG3	1:A:147:GLU:O	1.99	0.61
1:A:36:ILE:CG1	1:A:37:THR:N	2.62	0.61
1:A:103:ILE:C	1:A:103:ILE:CD1	2.58	0.61
1:A:117:LEU:CD2	1:B:117:LEU:HD23	2.28	0.61
1:B:257:PHE:O	1:B:261:GLU:CB	2.49	0.61
1:B:75:GLU:OE1	1:B:75:GLU:HA	2.01	0.60
1:A:18:LYS:CE	1:A:47:MET:SD	2.89	0.60
1:A:34:SER:O	1:A:38:HIS:CB	2.50	0.60
1:A:155:SER:O	1:A:156:ASN:C	2.40	0.60
1:A:228:GLY:CA	1:A:273:ILE:HD11	2.32	0.59
1:A:103:ILE:CG1	1:A:104:PHE:N	2.64	0.59
1:A:100:VAL:O	1:A:103:ILE:HG23	2.03	0.59
1:B:217:TYR:CE1	1:B:276:THR:HG21	2.38	0.59
1:B:243:ALA:O	1:B:247:THR:HB	2.02	0.59
1:A:122:VAL:HG13	2:B:333:HOH:O	2.02	0.58
1:B:275:PHE:CD1	1:B:279:ASN:ND2	2.71	0.58
1:A:250:ALA:O	1:A:254:GLN:HB2	2.02	0.58
1:A:160:SER:OG	1:A:163:VAL:HG23	2.03	0.58
1:A:76:LEU:HB2	1:A:120[A]:SER:OG	2.02	0.58
1:B:3:ILE:H	1:B:3:ILE:HD12	1.67	0.58
1:A:36:ILE:HD12	1:A:37:THR:CA	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:PHE:O	1:A:105:ASN:C	2.41	0.58
1:B:231:MET:O	1:B:259:VAL:CG1	2.51	0.58
1:B:290:GLU:HG3	2:B:350:HOH:O	2.04	0.57
1:A:135:LEU:C	1:A:135:LEU:HD23	2.25	0.57
1:B:243:ALA:CB	1:B:275:PHE:CE2	2.88	0.56
1:B:251:TRP:CD1	1:B:255:ILE:CD1	2.84	0.56
1:A:29:ASN:HB3	1:A:32:LEU:H	1.71	0.56
1:B:3:ILE:N	1:B:3:ILE:HD12	2.21	0.56
1:B:227:THR:CG2	1:B:228:GLY:N	2.68	0.56
1:A:36:ILE:HG13	1:A:37:THR:N	2.21	0.55
1:A:59:LEU:HD22	1:A:303:LEU:HD13	1.88	0.55
1:B:275:PHE:CD2	1:B:276:THR:N	2.74	0.55
1:B:159:PHE:CE2	1:B:244:LEU:HD13	2.41	0.55
1:A:18:LYS:HE2	1:A:47:MET:SD	2.47	0.55
1:A:18:LYS:HE3	1:A:47:MET:SD	2.47	0.55
1:A:163:VAL:O	1:A:167:VAL:HG23	2.06	0.55
1:A:252:ALA:O	1:A:254:GLN:N	2.40	0.54
1:B:265:THR:O	1:B:267:ASP:N	2.40	0.54
1:B:251:TRP:CG	1:B:252:ALA:N	2.75	0.53
1:A:164:TYR:CE2	1:A:238:LEU:HB2	2.43	0.53
1:B:32:LEU:O	1:B:34:ASP:N	2.42	0.53
1:B:227:THR:HG22	1:B:228:GLY:N	2.23	0.53
1:A:114:ASP:OD2	1:B:84:HIS:ND1	2.40	0.53
1:A:88:VAL:CG1	1:A:89:ASP:OD1	2.40	0.53
1:A:252:ALA:C	1:A:254:GLN:N	2.62	0.52
1:B:266:PRO:O	1:B:270:VAL:HG23	2.09	0.52
1:A:84:HIS:ND1	1:B:114:ASP:OD2	2.29	0.52
1:A:148:LEU:HD13	1:B:111:LEU:HB3	1.92	0.52
1:A:211:LYS:HD2	1:A:320:VAL:O	2.09	0.52
1:A:192:GLU:HG2	1:A:193:GLU:N	2.25	0.52
1:A:3:ILE:HD11	1:A:318:TYR:CE1	2.45	0.51
1:B:135:LEU:HD23	1:B:135:LEU:C	2.31	0.51
1:A:28:SER:N	1:B:145:GLU:OE1	2.43	0.51
1:B:243:ALA:HB3	1:B:275:PHE:CE2	2.46	0.51
1:A:252:ALA:C	1:A:254:GLN:H	2.13	0.51
1:B:217:TYR:HE1	1:B:276:THR:HG21	1.75	0.51
1:A:39:ILE:HD11	1:A:116:LEU:HD11	1.92	0.50
1:A:32:LEU:CD2	1:B:145:GLU:HG3	2.18	0.50
1:B:319:VAL:O	1:B:319:VAL:CG1	2.60	0.50
1:B:243:ALA:HB1	1:B:275:PHE:CE2	2.47	0.50
1:A:161:GLU:OE2	1:A:284:TYR:OH	2.30	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:278:ASP:OD1	1:B:278:ASP:O	2.30	0.50
1:A:130:TYR:CE2	1:B:130:TYR:CE2	2.99	0.49
1:B:3:ILE:HG23	1:B:7:LYS:HE3	1.94	0.49
1:B:150:GLN:O	1:B:153:ASN:HB3	2.11	0.49
1:B:267:ASP:O	1:B:270:VAL:CB	2.56	0.48
1:B:32:LEU:O	1:B:35:SER:N	2.36	0.48
1:A:149:LEU:O	1:A:153:ASN:ND2	2.45	0.48
1:B:212:ASP:OD1	1:B:322:ARG:NH2	2.46	0.48
1:B:234:GLY:HA2	1:B:259:VAL:HG12	1.95	0.48
1:B:211:LYS:CB	1:B:322:ARG:HH12	2.27	0.48
1:B:251:TRP:NE1	1:B:255:ILE:HD11	2.28	0.48
1:B:270:VAL:CG1	1:B:271:ARG:N	2.73	0.48
1:A:36:ILE:HD12	1:A:37:THR:HA	1.95	0.48
1:B:49:ARG:CD	1:B:175:LEU:CD1	2.85	0.48
1:B:265:THR:H	1:B:268:GLU:HG3	1.79	0.47
1:A:85:ASP:OD1	1:A:147:GLU:OE1	2.32	0.47
1:B:135:LEU:HD23	1:B:135:LEU:O	2.14	0.47
1:A:219:ASP:O	1:A:220:SER:C	2.52	0.47
1:A:8:THR:N	1:A:9:PRO:CD	2.76	0.47
1:B:229:ASN:ND2	1:B:232:LEU:HD23	2.29	0.47
1:A:196:PHE:CD1	1:A:196:PHE:C	2.88	0.47
1:A:89:ASP:N	1:A:89:ASP:OD1	2.47	0.47
1:B:49:ARG:HB2	1:B:50:PRO:HD3	1.96	0.47
1:A:106:ASN:ND2	1:A:106:ASN:N	2.58	0.47
1:B:189:VAL:HG23	1:B:193:GLU:HB2	1.96	0.47
1:A:76:LEU:HB3	1:A:120[A]:SER:HG	1.78	0.47
1:B:129:ASN:OD1	1:B:129:ASN:C	2.53	0.47
1:B:138:SER:O	1:B:141:GLN:HB2	2.14	0.46
1:B:206:ILE:O	1:B:210:ILE:HG13	2.15	0.46
1:A:36:ILE:CD1	1:A:36:ILE:O	2.48	0.46
1:B:269:ILE:O	1:B:273:ILE:CG1	2.48	0.46
1:B:12:ALA:HB3	2:B:343:HOH:O	2.15	0.46
1:B:319:VAL:O	1:B:319:VAL:HG12	2.15	0.46
1:B:231:MET:HA	1:B:259:VAL:HG11	1.97	0.46
1:A:283:GLU:CD	1:A:283:GLU:H	2.19	0.46
1:A:29:ASN:HD21	1:B:148:LEU:C	2.19	0.46
1:B:139:LEU:HD21	1:B:175:LEU:HD23	1.97	0.45
1:A:229:ASN:C	1:A:229:ASN:ND2	2.68	0.45
1:A:111:LEU:HD13	1:B:147:GLU:HB3	1.98	0.45
1:B:268:GLU:C	1:B:270:VAL:H	2.20	0.45
1:B:3:ILE:H	1:B:3:ILE:CD1	2.30	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:63:VAL:HG13	1:B:67:THR:HB	1.99	0.45
1:A:3:ILE:CD1	1:A:3:ILE:N	2.79	0.45
1:B:190:GLY:O	1:B:194:VAL:HG23	2.17	0.44
1:B:263:THR:OG1	1:B:263:THR:O	2.30	0.44
1:A:152:SER:O	1:A:153:ASN:C	2.56	0.44
1:B:229:ASN:O	1:B:232:LEU:HB3	2.18	0.44
1:B:135:LEU:HD21	1:B:178:ALA:HB1	2.00	0.44
1:B:234:GLY:HA2	1:B:259:VAL:O	2.18	0.43
1:A:64:THR:HB	1:A:65:PRO:HD2	2.00	0.43
1:B:118:ALA:O	1:B:122:VAL:HG23	2.17	0.43
1:A:78:HIS:HB2	1:A:175:LEU:HD21	2.00	0.43
1:B:287:ARG:O	1:B:291:GLN:HG3	2.17	0.43
1:B:192:GLU:CD	1:B:192:GLU:H	2.22	0.43
1:B:211:LYS:HG2	1:B:289:ILE:HD13	1.99	0.43
1:B:220:SER:O	1:B:221:LYS:C	2.57	0.43
1:A:210:ILE:HG22	1:A:289:ILE:HD11	1.99	0.43
1:A:159:PHE:CD1	1:A:159:PHE:N	2.85	0.43
1:B:244:LEU:HD23	1:B:275:PHE:CZ	2.53	0.43
1:A:165:PHE:HE2	1:A:210:ILE:HD11	1.84	0.43
1:B:242:TYR:CD2	1:B:242:TYR:C	2.93	0.42
1:B:84:HIS:CE1	1:B:114:ASP:OD1	2.72	0.42
1:B:8:THR:HB	1:B:9:PRO:HD3	2.01	0.42
1:B:32:LEU:C	1:B:34:ASP:N	2.73	0.42
1:A:135:LEU:HD23	1:A:136:VAL:N	2.34	0.42
1:A:60:TYR:HB2	1:A:187:VAL:HG21	2.02	0.42
1:A:130:TYR:CZ	1:B:130:TYR:CE2	3.07	0.42
1:A:64:THR:HB	1:A:65:PRO:CD	2.50	0.42
1:B:318:TYR:CD2	1:B:318:TYR:C	2.92	0.42
1:A:88:VAL:HG23	1:B:107:LYS:CB	2.50	0.42
1:B:251:TRP:O	1:B:255:ILE:HG13	2.20	0.42
1:A:255:ILE:O	1:A:259:VAL:HG23	2.20	0.42
1:A:247:THR:O	1:A:248:LYS:CB	2.67	0.42
1:B:5:LEU:HA	1:B:5:LEU:HD23	1.81	0.42
1:A:238:LEU:N	1:A:239:PRO:CD	2.83	0.41
1:B:148:LEU:O	1:B:152:SER:HB2	2.19	0.41
1:B:268:GLU:C	1:B:270:VAL:N	2.73	0.41
1:B:189:VAL:HG21	1:B:193:GLU:HB3	2.02	0.41
1:A:29:ASN:CB	1:A:32:LEU:HB2	2.49	0.41
1:B:211:LYS:HD2	1:B:320:VAL:O	2.20	0.41
1:B:244:LEU:HD11	1:B:256:ALA:HB2	2.01	0.41
1:A:189:VAL:HB	1:A:193:GLU:OE1	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:ARG:HB2	1:A:50:PRO:HD3	2.02	0.41
1:A:32:LEU:HD11	1:A:115:TYR:CE2	2.56	0.41
1:B:257:PHE:O	1:B:261:GLU:N	2.48	0.41
1:B:258:LYS:CB	1:B:263:THR:O	2.68	0.41
1:A:252:ALA:O	1:A:253:GLU:C	2.58	0.41
1:A:17:PHE:C	1:A:17:PHE:CD2	2.94	0.41
1:B:169:ARG:O	1:B:174:ALA:HB2	2.20	0.41
1:A:228:GLY:HA3	1:A:273:ILE:HD11	2.03	0.41
1:A:305:ASP:OD2	1:A:309:CYS:HB3	2.20	0.41
1:B:153:ASN:O	1:B:153:ASN:OD1	2.39	0.41
1:A:257:PHE:O	1:A:261:GLU:HB2	2.21	0.41
1:B:84:HIS:O	1:B:87:VAL:HG12	2.21	0.40
1:A:35:VAL:HG11	1:A:112:ALA:HB2	2.03	0.40
1:A:39:ILE:CD1	1:A:116:LEU:HD11	2.51	0.40
1:B:148:LEU:HD13	1:B:148:LEU: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	288/334 (86%)	277 (96%)	10 (4%)	1 (0%)	46	57
1	B	287/334 (86%)	267 (93%)	18 (6%)	2 (1%)	26	31
All	All	575/668 (86%)	544 (95%)	28 (5%)	3 (0%)	34	41

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	33	LEU
1	A	253	GLU
1	B	266	PRO

### 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	246/282 (87%)	234 (95%)	12 (5%)	31	41
1	B	235/282 (83%)	221 (94%)	14 (6%)	24	31
All	All	481/564 (85%)	455 (95%)	26 (5%)	27	36

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ILE
1	A	27	ASP
1	A	36	ILE
1	A	37	THR
1	A	89	ASP
1	A	103	ILE
1	A	147	GLU
1	A	153	ASN
1	A	196	PHE
1	A	211	LYS
1	A	229	ASN
1	A	265	THR
1	B	35	SER
1	B	39	HIS
1	B	152	SER
1	B	157	HIS
1	B	188	GLN
1	B	211	LYS
1	B	212	ASP
1	B	231	MET
1	B	247	THR
1	B	251	TRP
1	B	267	ASP
1	B	270	VAL
1	B	275	PHE
1	B	297	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	ASN
1	A	106	ASN
1	A	229	ASN
1	A	245	ASN
1	A	291	GLN
1	B	106	ASN
1	B	126	GLN
1	B	141	GLN
1	B	150	GLN
1	B	229	ASN
1	B	279	ASN
1	B	291	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	299/334 (89%)	0.48	27 (9%) 12 17	31, 59, 110, 132	0
1	B	297/334 (88%)	0.66	36 (12%) 6 9	31, 63, 122, 133	0
All	All	596/668 (89%)	0.57	63 (10%) 8 12	31, 60, 118, 133	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	31	ALA	8.0
1	A	154	VAL	8.0
1	A	32	LEU	7.4
1	B	280	GLY	6.8
1	A	100	VAL	6.8
1	A	101	ASN	6.6
1	B	154	VAL	6.4
1	A	40	ARG	6.3
1	A	151	LEU	6.1
1	A	256	ALA	5.6
1	A	99	SER	5.5
1	B	103	ILE	5.1
1	B	281	GLY	5.0
1	A	257	PHE	4.9
1	B	39	HIS	4.9
1	A	36	ILE	4.6
1	A	111	LEU	4.6
1	A	103	ILE	4.6
1	B	275	PHE	4.5
1	A	102	ALA	4.4
1	B	282	ILE	4.3
1	A	149	LEU	4.0
1	A	104	PHE	3.9
1	B	269	ILE	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	33	LEU	3.9
1	B	278	ASP	3.8
1	B	155	SER	3.8
1	B	246	THR	3.7
1	B	219	ASP	3.6
1	B	271	ARG	3.5
1	A	108	VAL	3.5
1	A	30	ALA	3.5
1	A	46	LYS	3.4
1	B	283	GLU	3.4
1	B	0	LEU	3.3
1	B	156	ASN	3.3
1	B	322	ARG	3.2
1	B	149	LEU	3.2
1	B	32	LEU	3.2
1	A	155	SER	3.1
1	B	232	LEU	3.0
1	B	279	ASN	3.0
1	A	156	ASN	2.9
1	B	254	GLN	2.8
1	B	247	THR	2.7
1	B	150	GLN	2.6
1	A	38	HIS	2.6
1	B	104	PHE	2.5
1	A	269	ILE	2.5
1	B	36	VAL	2.5
1	B	287	ARG	2.4
1	A	159	PHE	2.4
1	A	250	ALA	2.3
1	B	236	LEU	2.3
1	B	272	LEU	2.3
1	B	274	GLU	2.3
1	A	272	LEU	2.2
1	B	102	ALA	2.2
1	B	130	TYR	2.1
1	B	262	GLY	2.1
1	B	153	ASN	2.1
1	A	35	VAL	2.0
1	A	251	TRP	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.