



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 04:58 am GMT

PDB ID : 1WXQ
Title : Crystal Structure of GTP binding protein from *Pyrococcus horikoshii* OT3
Authors : Lokanath, N.K.; Kunishima, N.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2005-01-29
Resolution : 2.60 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

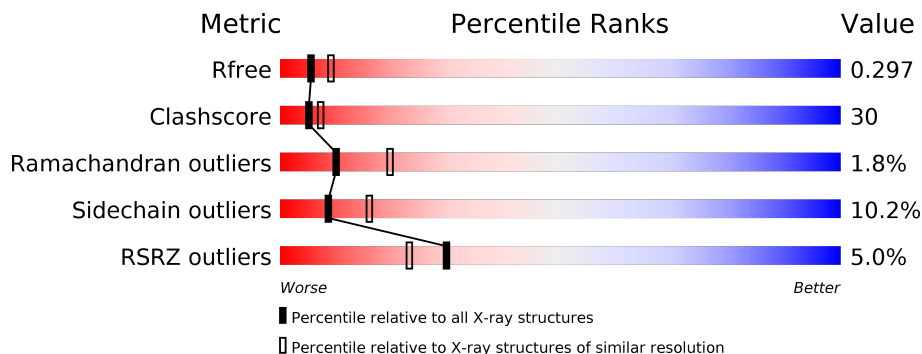
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	100719	2542 (2.60-2.60)
Clashscore	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)
RSRZ outliers	101464	2550 (2.60-2.60)

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	397	<div> <div>4%</div> <div>48%</div> <div>31%</div> <div>7%</div> <div>13%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2807 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 GTP-binding protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	344	Total	C	N	O	S	Se	0	0	0
			2718	1744	462	503	2	7			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP O58261
A	75	MSE	MET	MODIFIED RESIDUE	UNP O58261
A	101	MSE	MET	MODIFIED RESIDUE	UNP O58261
A	186	MSE	MET	MODIFIED RESIDUE	UNP O58261
A	217	MSE	MET	MODIFIED RESIDUE	UNP O58261
A	281	MSE	MET	MODIFIED RESIDUE	UNP O58261
A	290	MSE	MET	MODIFIED RESIDUE	UNP O58261
A	345	MSE	MET	MODIFIED RESIDUE	UNP O58261

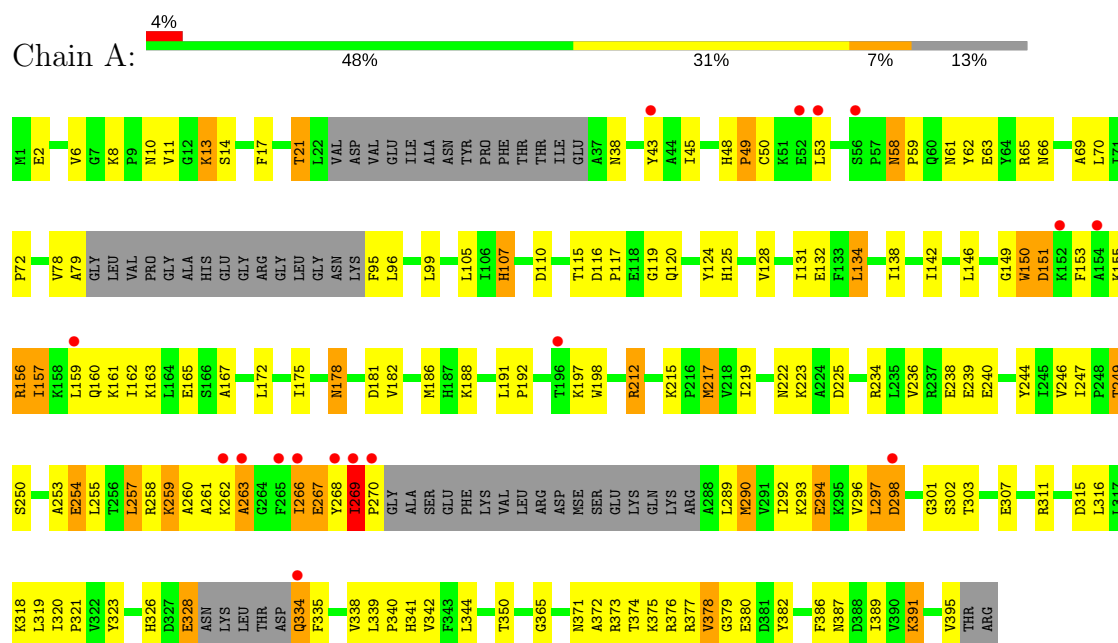
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	89	Total	O	0	0
			89	89		

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: GTP-binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	74.96Å 143.00Å 78.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.60 19.71 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.6 (10.00-2.60) 97.6 (19.71-2.40)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.16 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.1.27	Depositor
R, R_{free}	0.266 , 0.292 0.271 , 0.297	Depositor DCC
R_{free} test set	1322 reflections (11.30%)	DCC
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.281	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 43.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	2807	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.31% 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	0.54	1/2765 (0.0%)	0.84	2/3729 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	150	TRP	NE1-CE2	8.78	1.49	1.37

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	269	ILE	CG1-CB-CG2	-5.92	98.37	111.40
1	A	391	LYS	CB-CG-CD	5.57	126.08	111.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	261	ALA	Peptide

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	2718	0	2751	162	0
2	A	89	0	0	10	0
All	All	2807	0	2751	162	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 30.

All (162) 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:53:LEU:HD23	1:A:53:LEU:O	1.27	1.28
1:A:17:PHE:O	1:A:21:THR:OG1	1.71	1.05
1:A:253:ALA:HA	1:A:293:LYS:HZ2	1.22	1.05
1:A:269:ILE:HD12	2:A:432:HOH:O	1.62	1.00
1:A:159:LEU:O	1:A:159:LEU:HD23	1.63	0.98
1:A:253:ALA:HA	1:A:293:LYS:NZ	1.88	0.88
1:A:269:ILE:HB	1:A:270:PRO:HD3	1.57	0.87
1:A:269:ILE:CB	1:A:270:PRO:HD3	2.04	0.87
1:A:53:LEU:HD23	1:A:53:LEU:C	1.96	0.86
1:A:334:GLN:HG2	1:A:335:PHE:CE1	2.11	0.85
1:A:17:PHE:CZ	1:A:21:THR:HG21	2.11	0.85
1:A:321:PRO:HD3	1:A:386:PHE:CE2	2.12	0.85
1:A:159:LEU:C	1:A:159:LEU:HD23	1.97	0.85
1:A:53:LEU:CD2	1:A:53:LEU:O	2.22	0.82
1:A:107:HIS:CD2	1:A:134:LEU:HD12	2.15	0.81
1:A:128:VAL:HG11	1:A:238:GLU:OE2	1.79	0.81
1:A:66:ASN:HB3	2:A:420:HOH:O	1.81	0.81
1:A:61:ASN:ND2	1:A:341:HIS:HB3	1.97	0.80
1:A:153:PHE:O	1:A:157:ILE:HD12	1.82	0.79
1:A:222:ASN:OD1	1:A:249:THR:HG22	1.82	0.79
1:A:268:TYR:O	1:A:269:ILE:HG13	1.83	0.79
1:A:254:GLU:O	1:A:258:ARG:HB2	1.82	0.79
1:A:45:ILE:HG12	1:A:70:LEU:HG	1.65	0.77
1:A:269:ILE:CG2	1:A:270:PRO:HD3	2.15	0.77
1:A:131:ILE:HD13	1:A:219:ILE:HD13	1.68	0.76
1:A:6:VAL:HG21	1:A:134:LEU:HD11	1.70	0.73
1:A:138:ILE:O	1:A:142:ILE:HG13	1.89	0.73
1:A:339:LEU:N	1:A:340:PRO:HD3	2.04	0.73
1:A:293:LYS:O	1:A:297:LEU:HB2	1.89	0.72

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:ASN:HD21	1:A:342:VAL:H	1.35	0.72
1:A:157:ILE:HG23	1:A:162:ILE:HB	1.71	0.72
1:A:292:ILE:HG13	1:A:296:VAL:CG2	2.20	0.71
1:A:269:ILE:CB	1:A:270:PRO:CD	2.68	0.71
1:A:269:ILE:HB	1:A:270:PRO:CD	2.20	0.70
1:A:146:LEU:HD21	1:A:172:LEU:HD11	1.75	0.69
1:A:257:LEU:O	1:A:266:ILE:HG13	1.92	0.69
1:A:263:ALA:HB2	2:A:479:HOH:O	1.92	0.68
1:A:159:LEU:CD2	1:A:159:LEU:C	2.64	0.66
1:A:53:LEU:CD2	1:A:53:LEU:C	2.64	0.66
1:A:61:ASN:HD21	1:A:341:HIS:HB3	1.60	0.66
1:A:290:MSE:O	1:A:290:MSE:HG2	1.96	0.65
1:A:59:PRO:HB3	1:A:62:TYR:CZ	2.31	0.65
1:A:378:VAL:HG13	1:A:379:GLY:O	1.96	0.65
1:A:269:ILE:HG22	1:A:270:PRO:N	2.11	0.65
1:A:128:VAL:O	1:A:128:VAL:HG12	1.96	0.64
1:A:269:ILE:HG22	1:A:270:PRO:CD	2.28	0.64
1:A:318:LYS:HD3	2:A:456:HOH:O	1.98	0.63
1:A:13:LYS:NZ	1:A:79:ALA:HB2	2.13	0.63
1:A:292:ILE:HG13	1:A:296:VAL:HG23	1.79	0.63
1:A:376:ARG:HG2	1:A:376:ARG:HH11	1.63	0.62
1:A:250:SER:HB2	1:A:297:LEU:HD21	1.81	0.62
1:A:253:ALA:O	1:A:257:LEU:HD12	2.00	0.62
1:A:217:MSE:HG3	1:A:244:TYR:HE1	1.65	0.61
1:A:117:PRO:HG3	1:A:125:HIS:NE2	2.16	0.61
1:A:236:VAL:O	1:A:240:GLU:HG3	2.01	0.61
1:A:254:GLU:HG3	1:A:303:THR:OG1	2.01	0.60
1:A:38:ASN:HB3	2:A:416:HOH:O	2.00	0.60
1:A:178:ASN:HB2	1:A:181:ASP:OD2	2.01	0.60
1:A:188:LYS:NZ	1:A:212:ARG:HH22	1.99	0.60
1:A:61:ASN:OD1	1:A:62:TYR:HD1	1.85	0.60
1:A:156:ARG:C	1:A:157:ILE:HG13	2.22	0.59
1:A:373:ARG:NH1	1:A:387:ASN:O	2.35	0.59
1:A:61:ASN:HD21	1:A:342:VAL:N	2.00	0.59
1:A:192:PRO:HD2	1:A:198:TRP:NE1	2.17	0.58
1:A:255:LEU:O	1:A:259:LYS:HB2	2.03	0.58
1:A:266:ILE:HB	1:A:268:TYR:CD1	2.37	0.58
1:A:146:LEU:O	1:A:150:TRP:HB2	2.04	0.57
1:A:156:ARG:O	1:A:157:ILE:HG13	2.05	0.57
1:A:63:GLU:OE2	1:A:65:ARG:HD2	2.05	0.57
1:A:321:PRO:CD	1:A:386:PHE:CE2	2.87	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:292:ILE:HG13	1:A:296:VAL:HG21	1.87	0.57
1:A:269:ILE:CG2	1:A:270:PRO:CD	2.84	0.56
1:A:13:LYS:HZ2	1:A:79:ALA:HB2	1.70	0.56
1:A:292:ILE:O	1:A:296:VAL:N	2.38	0.55
1:A:188:LYS:HZ1	1:A:212:ARG:HH12	1.54	0.55
1:A:188:LYS:NZ	1:A:212:ARG:HH12	2.04	0.55
1:A:159:LEU:O	1:A:161:LYS:HG2	2.07	0.54
1:A:296:VAL:HG12	2:A:453:HOH:O	2.07	0.54
1:A:328:GLU:OE1	1:A:328:GLU:N	2.40	0.54
1:A:175:ILE:N	1:A:175:ILE:HD12	2.22	0.54
1:A:267:GLU:H	1:A:268:TYR:HA	1.71	0.54
1:A:163:LYS:HG3	1:A:165:GLU:HG2	1.90	0.54
1:A:128:VAL:CG1	1:A:128:VAL:O	2.55	0.53
1:A:99:LEU:O	1:A:215:LYS:HE3	2.08	0.53
1:A:315:ASP:O	1:A:318:LYS:HG2	2.10	0.52
1:A:162:ILE:CG2	1:A:163:LYS:N	2.73	0.52
1:A:146:LEU:HD21	1:A:172:LEU:CD1	2.39	0.52
1:A:266:ILE:HD13	1:A:266:ILE:H	1.75	0.51
1:A:307:GLU:HG2	2:A:406:HOH:O	2.10	0.51
1:A:116:ASP:HB2	1:A:120:GLN:O	2.10	0.51
1:A:292:ILE:O	1:A:296:VAL:HG23	2.11	0.51
1:A:258:ARG:C	1:A:260:ALA:H	2.14	0.51
1:A:128:VAL:HG13	1:A:239:GLU:OE2	2.09	0.51
1:A:294:GLU:O	1:A:298:ASP:HB2	2.10	0.51
1:A:365:GLY:CA	1:A:395:VAL:HG12	2.41	0.51
1:A:378:VAL:HG21	1:A:382:TYR:CD2	2.46	0.50
1:A:61:ASN:ND2	1:A:342:VAL:H	2.06	0.50
1:A:269:ILE:HG22	1:A:270:PRO:HD3	1.84	0.50
1:A:301:GLY:O	1:A:302:SER:HB3	2.11	0.50
1:A:371:ASN:OD1	1:A:373:ARG:HB2	2.12	0.50
1:A:186:MSE:HA	1:A:191:LEU:HD12	1.94	0.50
1:A:99:LEU:HD12	1:A:99:LEU:C	2.32	0.50
1:A:149:GLY:O	1:A:151:ASP:N	2.44	0.49
1:A:219:ILE:HB	1:A:246:VAL:HG22	1.93	0.49
1:A:105:LEU:HB2	1:A:217:MSE:HB3	1.93	0.49
1:A:378:VAL:HG13	1:A:379:GLY:N	2.28	0.49
1:A:62:TYR:N	1:A:62:TYR:CD1	2.82	0.48
1:A:266:ILE:HD13	1:A:266:ILE:O	2.13	0.48
1:A:58:ASN:OD1	1:A:338:VAL:HG12	2.14	0.48
1:A:372:ALA:HB3	1:A:389:ILE:HB	1.96	0.48
1:A:6:VAL:HG12	1:A:78:VAL:CG1	2.43	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:ILE:H	1:A:266:ILE:CD1	2.26	0.48
1:A:267:GLU:N	1:A:268:TYR:HA	2.29	0.47
1:A:8:LYS:HG2	2:A:408:HOH:O	2.14	0.47
1:A:320:ILE:HG22	1:A:386:PHE:HD2	1.79	0.47
1:A:124:TYR:OH	1:A:234:ARG:NH2	2.47	0.47
1:A:323:TYR:CE2	1:A:342:VAL:HB	2.50	0.47
1:A:319:LEU:HD13	1:A:344:LEU:HD13	1.95	0.47
1:A:49:PRO:HG3	1:A:323:TYR:CE1	2.50	0.46
1:A:6:VAL:HG12	1:A:78:VAL:HB	1.96	0.46
1:A:13:LYS:HG3	1:A:14:SER:N	2.22	0.46
1:A:339:LEU:N	1:A:340:PRO:CD	2.77	0.46
1:A:165:GLU:HA	1:A:182:VAL:HG11	1.97	0.45
1:A:11:VAL:HG12	1:A:110:ASP:HB2	1.99	0.45
1:A:328:GLU:HG3	1:A:391:LYS:HB2	1.98	0.45
1:A:175:ILE:HG22	1:A:175:ILE:O	2.17	0.45
1:A:376:ARG:HG2	1:A:376:ARG:NH1	2.31	0.45
1:A:374:THR:O	1:A:375:LYS:HB2	2.17	0.45
1:A:334:GLN:HG2	1:A:335:PHE:CD1	2.49	0.45
1:A:386:PHE:O	1:A:387:ASN:HB2	2.16	0.45
1:A:156:ARG:HB3	1:A:156:ARG:HE	1.58	0.44
1:A:326:HIS:O	1:A:391:LYS:HD3	2.16	0.44
1:A:155:LYS:C	1:A:157:ILE:H	2.21	0.44
1:A:192:PRO:HD2	1:A:198:TRP:CD1	2.52	0.44
1:A:250:SER:CB	1:A:297:LEU:HD21	2.48	0.44
1:A:290:MSE:HB3	2:A:460:HOH:O	2.18	0.44
1:A:247:ILE:HD11	1:A:311:ARG:HG3	2.00	0.44
1:A:128:VAL:HG21	1:A:238:GLU:CD	2.39	0.43
1:A:117:PRO:HG3	1:A:125:HIS:CE1	2.53	0.43
1:A:48:HIS:O	1:A:50:CYS:N	2.51	0.43
1:A:110:ASP:OD2	1:A:115:THR:HG23	2.17	0.43
1:A:212:ARG:HB3	1:A:212:ARG:HE	1.54	0.43
1:A:61:ASN:ND2	1:A:341:HIS:CB	2.76	0.43
1:A:10:ASN:ND2	1:A:119:GLY:O	2.36	0.43
1:A:162:ILE:HG22	1:A:163:LYS:N	2.32	0.43
1:A:225:ASP:N	1:A:225:ASP:OD1	2.51	0.42
1:A:311:ARG:O	1:A:315:ASP:HB3	2.20	0.42
1:A:165:GLU:CD	1:A:165:GLU:H	2.21	0.42
1:A:150:TRP:O	1:A:150:TRP:CG	2.73	0.42
1:A:45:ILE:HA	1:A:69:ALA:O	2.20	0.42
1:A:48:HIS:CG	1:A:49:PRO:HD2	2.54	0.41
1:A:350:THR:HG22	1:A:382:TYR:O	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:371:ASN:C	1:A:373:ARG:H	2.24	0.41
1:A:380:GLU:O	1:A:380:GLU:HG3	2.19	0.41
1:A:43:TYR:CE1	1:A:72:PRO:HB3	2.55	0.41
1:A:307:GLU:CG	2:A:406:HOH:O	2.67	0.41
1:A:376:ARG:HG2	1:A:377:ARG:N	2.36	0.41
1:A:222:ASN:O	1:A:223:LYS:HB2	2.21	0.41
1:A:266:ILE:N	1:A:266:ILE:CD1	2.84	0.41
1:A:255:LEU:HG	1:A:259:LYS:HE2	2.03	0.41
1:A:162:ILE:HG21	1:A:167:ALA:HB2	2.03	0.40
1:A:268:TYR:HE2	1:A:303:THR:HG23	1.85	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	334/397 (84%)	294 (88%)	34 (10%)	6 (2%)	10	19

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	269	ILE
1	A	259	LYS
1	A	263	ALA
1	A	262	LYS
1	A	157	ILE
1	A	49	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	293/329 (89%)	263 (90%)	30 (10%)	8 16

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLU
1	A	13	LYS
1	A	21	THR
1	A	58	ASN
1	A	95	PHE
1	A	96	LEU
1	A	107	HIS
1	A	132	GLU
1	A	134	LEU
1	A	151	ASP
1	A	156	ARG
1	A	160	GLN
1	A	178	ASN
1	A	197	LYS
1	A	212	ARG
1	A	217	MSE
1	A	249	THR
1	A	254	GLU
1	A	257	LEU
1	A	266	ILE
1	A	267	GLU
1	A	289	LEU
1	A	290	MSE
1	A	294	GLU
1	A	297	LEU
1	A	298	ASP
1	A	316	LEU
1	A	328	GLU
1	A	334	GLN
1	A	378	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	107	HIS
1	A	120	GLN
1	A	160	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 [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	337/397 (84%)	0.23	17 (5%) 30 23	18, 37, 61, 72	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	266	ILE	7.5
1	A	159	LEU	6.5
1	A	53	LEU	5.5
1	A	262	LYS	4.7
1	A	269	ILE	4.2
1	A	268	TYR	4.0
1	A	52	GLU	3.7
1	A	56	SER	2.9
1	A	152	LYS	2.8
1	A	298	ASP	2.8
1	A	265	PHE	2.6
1	A	43	TYR	2.5
1	A	196	THR	2.4
1	A	154	ALA	2.2
1	A	263	ALA	2.1
1	A	270	PRO	2.1
1	A	334	GLN	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.