



# Full wwPDB X-ray Structure Validation Report

Aug 22, 2023 – 03:13 AM EDT

PDB ID : 2PIG  
Title : Crystal structure of ydcK from Salmonella cholerae at 2.38 Å resolution.  
Northeast Structural Genomics target SCR6  
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Conover, K.; Ma, L.-C.; Xiao, R.; Acton, T.B.; Montelione, G.T.; Hunt, J.F.;  
Tong, L.; Northeast Structural Genomics Consortium (NESG)  
Deposited on : 2007-04-13  
Resolution : 2.38 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtrriage (Phenix) : 1.13  
EDS : 2.35  
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) : 2.35

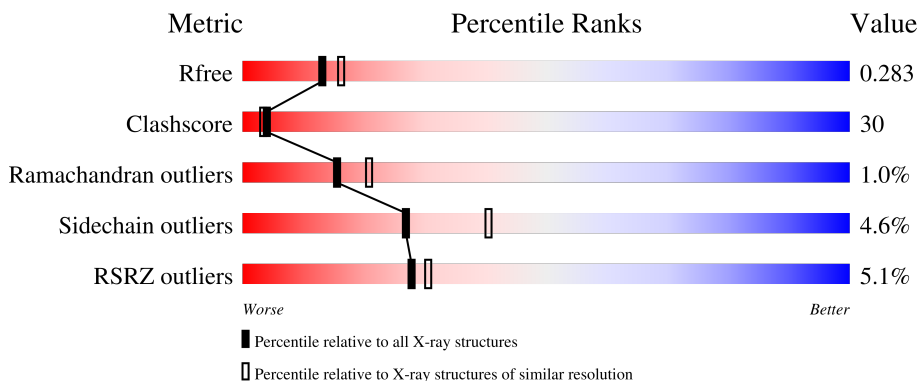
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.38 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	5509 (2.40-2.36)
Clashscore	141614	6082 (2.40-2.36)
Ramachandran outliers	138981	5973 (2.40-2.36)
Sidechain outliers	138945	5975 (2.40-2.36)
RSRZ outliers	127900	5397 (2.40-2.36)

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 of 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	334	
1	B	334	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5024 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 Putative transferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	310	2407	1495	437	467	7	1	0	0	0
1	B	319	2479	1544	450	477	7	1	0	0	0

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP Q5PHU6
A	67	MSE	MET	modified residue	UNP Q5PHU6
A	327	LEU	-	cloning artifact	UNP Q5PHU6
A	328	GLU	-	cloning artifact	UNP Q5PHU6
A	329	HIS	-	cloning artifact	UNP Q5PHU6
A	330	HIS	-	cloning artifact	UNP Q5PHU6
A	331	HIS	-	cloning artifact	UNP Q5PHU6
A	332	HIS	-	cloning artifact	UNP Q5PHU6
A	333	HIS	-	cloning artifact	UNP Q5PHU6
A	334	HIS	-	cloning artifact	UNP Q5PHU6
B	1	MSE	MET	modified residue	UNP Q5PHU6
B	67	MSE	MET	modified residue	UNP Q5PHU6
B	327	LEU	-	cloning artifact	UNP Q5PHU6
B	328	GLU	-	cloning artifact	UNP Q5PHU6
B	329	HIS	-	cloning artifact	UNP Q5PHU6
B	330	HIS	-	cloning artifact	UNP Q5PHU6
B	331	HIS	-	cloning artifact	UNP Q5PHU6
B	332	HIS	-	cloning artifact	UNP Q5PHU6
B	333	HIS	-	cloning artifact	UNP Q5PHU6
B	334	HIS	-	cloning artifact	UNP Q5PHU6

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total 2	Zn 2	0	0
2	B	1	Total 1	Zn 1	0	0

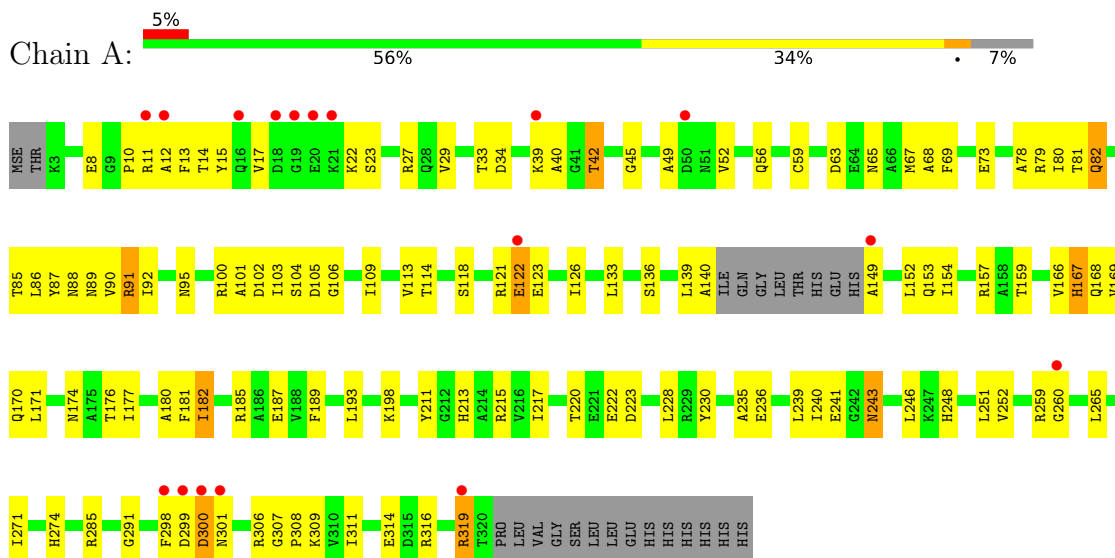
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	67	Total 67	O 67	0	0
3	B	68	Total 68	O 68	0	0

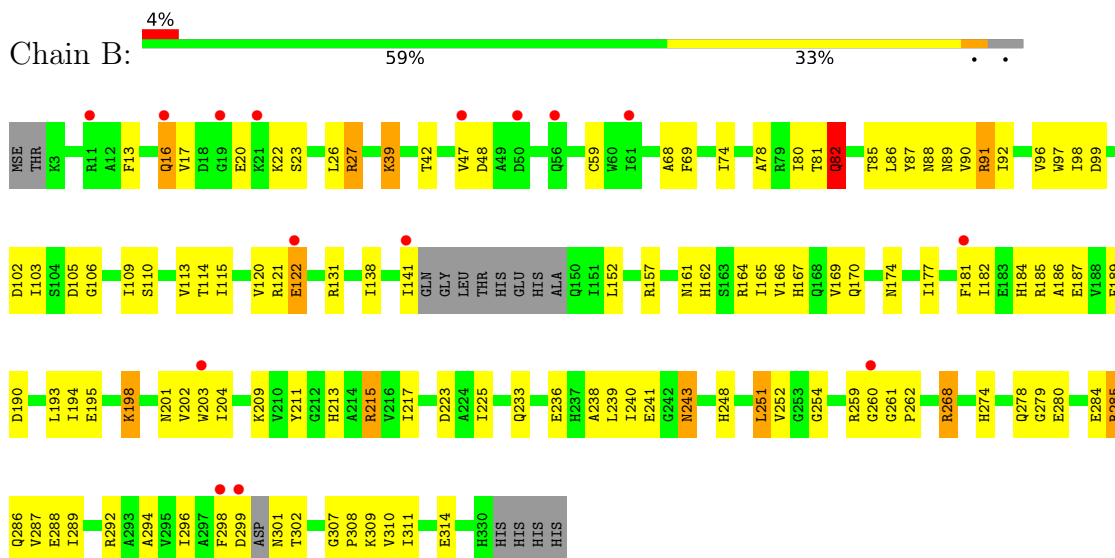
### 3 Residue-property plots i

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Putative transferase



- Molecule 1: Putative transferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	103.60Å 103.60Å 160.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.38 19.79 – 2.30	Depositor EDS
% Data completeness (in resolution range)	89.8 (20.00-2.38) 95.5 (19.79-2.30)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.33 (at 2.30Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.220 , 0.269 0.233 , 0.283	Depositor DCC
$R_{free}$ test set	3974 reflections (10.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.4	Xtrriage
Anisotropy	0.141	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 42.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	5024	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.49% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section:  
ZN

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	0.31	0/2444	0.65	0/3315
1	B	0.31	0/2518	0.63	0/3415
All	All	0.31	0/4962	0.64	0/6730

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	2407	0	2347	137	0
1	B	2479	0	2426	163	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
3	A	67	0	0	16	0
3	B	68	0	0	11	0
All	All	5024	0	4773	292	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 (292) 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:287:VAL:HG12	1:B:289:ILE:HD11	1.21	1.16
1:B:141:ILE:HD12	3:B:3108:HOH:O	1.43	1.14
1:A:73:GLU:HG3	1:A:91:ARG:HH12	1.09	1.13
1:A:309:LYS:HE3	1:A:311:ILE:HD11	1.13	1.09
1:B:90:VAL:CG1	1:B:92:ILE:HD11	1.82	1.08
1:B:287:VAL:HG12	1:B:289:ILE:CD1	1.85	1.06
1:B:285:ARG:HG3	1:B:285:ARG:HH11	0.88	1.04
1:A:39:LYS:O	1:A:42:THR:HG23	1.56	1.04
1:B:91:ARG:C	1:B:92:ILE:HD12	1.79	1.02
1:B:287:VAL:CG1	1:B:289:ILE:HD11	1.90	1.02
1:B:285:ARG:HH11	1:B:285:ARG:CG	1.74	1.01
1:B:285:ARG:HG3	1:B:285:ARG:NH1	1.69	0.98
1:A:309:LYS:HE3	1:A:311:ILE:CD1	1.93	0.98
1:B:294:ALA:HB1	1:B:296:ILE:HD11	1.45	0.97
1:A:259:ARG:C	3:A:3010:HOH:O	2.02	0.97
1:B:90:VAL:HG12	1:B:92:ILE:CD1	1.95	0.96
1:B:170:GLN:HB2	1:B:187:GLU:HG2	1.46	0.95
1:A:39:LYS:O	1:A:42:THR:CG2	2.17	0.93
1:A:73:GLU:HG3	1:A:91:ARG:NH1	1.84	0.92
1:A:149:ALA:N	3:A:3135:HOH:O	2.02	0.91
1:B:141:ILE:HG22	1:B:167:HIS:HA	1.53	0.91
1:B:298:PHE:O	1:B:301:ASN:HB2	1.72	0.90
1:B:85:THR:HB	1:B:102:ASP:OD1	1.76	0.83
1:A:306:ARG:NE	3:A:3125:HOH:O	2.11	0.83
1:B:288:GLU:C	1:B:289:ILE:HD12	2.02	0.81
1:B:169:VAL:HG21	1:B:182:ILE:HG22	1.62	0.80
1:A:260:GLY:N	3:A:3010:HOH:O	2.12	0.80
1:B:236:GLU:OE1	1:B:254:GLY:N	2.11	0.80
1:A:177:ILE:HD12	1:A:177:ILE:N	1.97	0.79
1:B:90:VAL:HG12	1:B:92:ILE:HD11	1.53	0.79
1:B:92:ILE:HD12	1:B:92:ILE:N	1.97	0.78
1:A:154:ILE:HD12	1:A:171:LEU:HD12	1.66	0.78
1:A:102:ASP:C	1:A:103:ILE:HD12	2.05	0.77
1:B:90:VAL:CG1	1:B:92:ILE:CD1	2.56	0.76
1:B:285:ARG:CG	1:B:285:ARG:NH1	2.39	0.76
1:B:82:GLN:HE21	1:B:82:GLN:N	1.84	0.76
1:B:48:ASP:OD2	1:B:69:PHE:HB2	1.87	0.75
1:A:8:GLU:HA	3:A:3115:HOH:O	1.87	0.74
1:A:169:VAL:HG21	1:A:182:ILE:HG22	1.69	0.73
1:B:164:ARG:HB3	1:B:181:PHE:HD1	1.53	0.73
1:B:202:VAL:HG12	1:B:204:ILE:CD1	2.18	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:ILE:HA	3:B:3108:HOH:O	1.88	0.73
1:B:169:VAL:CG2	1:B:182:ILE:HG22	2.18	0.73
1:B:27:ARG:HH11	1:B:27:ARG:HG3	1.54	0.73
1:B:202:VAL:HG12	1:B:204:ILE:HD11	1.70	0.73
1:B:289:ILE:HD12	1:B:289:ILE:N	2.03	0.72
1:A:309:LYS:CE	1:A:311:ILE:HD11	2.07	0.71
1:B:203:TRP:C	1:B:204:ILE:HD12	2.10	0.70
1:B:27:ARG:HG3	1:B:27:ARG:NH1	2.06	0.70
3:A:3038:HOH:O	1:B:131:ARG:HD3	1.90	0.70
1:A:170:GLN:HB2	1:A:187:GLU:HG2	1.74	0.69
1:B:39:LYS:HB2	1:B:39:LYS:NZ	2.08	0.69
1:B:152:LEU:HD23	1:B:152:LEU:C	2.12	0.68
1:A:85:THR:HG22	1:A:87:TYR:CE1	2.27	0.68
1:B:166:VAL:HG12	1:B:167:HIS:N	2.08	0.68
1:B:82:GLN:N	1:B:82:GLN:NE2	2.40	0.68
1:B:165:ILE:N	1:B:165:ILE:HD12	2.09	0.68
1:B:274:HIS:CE1	3:B:3071:HOH:O	2.47	0.68
1:B:22:LYS:NZ	1:B:22:LYS:HB3	2.09	0.67
1:A:251:LEU:HD23	1:A:251:LEU:C	2.15	0.67
1:B:164:ARG:HB3	1:B:181:PHE:CD1	2.29	0.67
1:B:90:VAL:HG11	1:B:92:ILE:HD11	1.75	0.67
1:A:29:VAL:HG22	1:A:45:GLY:O	1.95	0.67
1:A:215:ARG:HD2	1:B:193:LEU:HD13	1.77	0.66
1:A:85:THR:HG22	1:A:87:TYR:CD1	2.31	0.66
1:B:141:ILE:HG23	1:B:141:ILE:O	1.96	0.66
1:B:239:LEU:C	1:B:239:LEU:HD23	2.16	0.66
1:B:274:HIS:HE1	3:B:3071:HOH:O	1.78	0.65
1:A:159:THR:HG22	3:B:3061:HOH:O	1.96	0.64
1:A:85:THR:C	1:A:86:LEU:HD12	2.17	0.64
1:A:157:ARG:NE	1:A:174:ASN:OD1	2.28	0.64
1:A:239:LEU:HD23	1:A:239:LEU:C	2.18	0.64
1:A:176:THR:C	1:A:177:ILE:HD12	2.18	0.64
1:B:311:ILE:N	1:B:311:ILE:HD12	2.13	0.64
1:B:193:LEU:HB3	1:B:215:ARG:HG3	1.79	0.64
1:B:86:LEU:N	1:B:86:LEU:HD12	2.12	0.64
1:A:80:ILE:N	1:A:80:ILE:HD12	2.13	0.63
1:B:204:ILE:HD12	1:B:204:ILE:N	2.14	0.63
1:A:86:LEU:HD12	1:A:86:LEU:N	2.14	0.62
1:A:260:GLY:C	3:A:3010:HOH:O	2.37	0.62
1:A:228:LEU:HD22	1:A:246:LEU:HD22	1.81	0.61
1:B:85:THR:HG22	1:B:87:TYR:CE2	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:PHE:C	1:A:182:ILE:HD12	2.22	0.60
1:A:211:TYR:CZ	1:A:235:ALA:HB2	2.35	0.60
1:B:177:ILE:N	1:B:177:ILE:HD12	2.16	0.60
1:A:121:ARG:O	1:A:122:GLU:HB2	2.00	0.60
1:B:166:VAL:HG12	1:B:167:HIS:CD2	2.37	0.60
1:B:309:LYS:HG2	1:B:311:ILE:HD11	1.84	0.60
1:B:182:ILE:HG23	1:B:186:ALA:HB3	1.84	0.59
1:A:85:THR:CG2	1:A:87:TYR:CE1	2.85	0.59
1:B:157:ARG:HH21	1:B:174:ASN:ND2	2.01	0.59
1:A:79:ARG:C	1:A:80:ILE:HD12	2.23	0.59
1:A:243:ASN:C	1:A:243:ASN:HD22	2.06	0.59
1:A:177:ILE:N	1:A:177:ILE:CD1	2.66	0.58
1:B:22:LYS:HB3	1:B:22:LYS:HZ2	1.68	0.58
1:A:301:ASN:HB2	3:A:3126:HOH:O	2.02	0.58
1:A:101:ALA:HB1	1:A:118:SER:O	2.04	0.58
1:B:121:ARG:O	1:B:122:GLU:HB2	2.03	0.58
1:B:285:ARG:HH12	1:B:308:PRO:HD3	1.69	0.57
1:B:294:ALA:CB	1:B:296:ILE:HD11	2.28	0.57
1:B:309:LYS:HG2	1:B:311:ILE:CD1	2.34	0.57
1:B:182:ILE:CG2	1:B:186:ALA:HB3	2.34	0.57
1:B:39:LYS:HB2	1:B:39:LYS:HZ3	1.68	0.57
1:B:296:ILE:HD12	1:B:296:ILE:N	2.18	0.57
1:A:59:CYS:HA	1:A:78:ALA:O	2.05	0.57
1:A:8:GLU:CB	3:A:3115:HOH:O	2.52	0.57
1:A:100:ARG:HD2	3:A:3029:HOH:O	2.04	0.57
1:A:169:VAL:CG2	1:A:182:ILE:HG22	2.33	0.56
1:B:141:ILE:CG2	1:B:167:HIS:HA	2.29	0.56
1:A:230:TYR:CE2	1:A:248:HIS:CE1	2.94	0.56
1:A:220:THR:OG1	1:B:198:LYS:HE2	2.05	0.56
1:A:298:PHE:O	1:A:300:ASP:N	2.38	0.56
1:B:85:THR:CG2	1:B:87:TYR:CE2	2.89	0.56
1:B:202:VAL:CG1	1:B:204:ILE:HD11	2.35	0.56
1:B:289:ILE:CD1	1:B:289:ILE:N	2.69	0.56
1:A:182:ILE:HD12	1:A:182:ILE:N	2.22	0.55
1:B:80:ILE:HD13	1:B:98:ILE:HB	1.87	0.55
1:A:14:THR:HA	1:A:22:LYS:O	2.07	0.55
1:B:85:THR:HG22	1:B:87:TYR:CD2	2.41	0.55
1:B:86:LEU:HA	1:B:103:ILE:O	2.06	0.55
1:B:261:GLY:HA2	1:B:262:PRO:C	2.26	0.55
1:A:90:VAL:C	1:A:91:ARG:HD2	2.28	0.55
1:B:166:VAL:CG1	1:B:167:HIS:N	2.70	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:251:LEU:C	1:B:251:LEU:HD23	2.27	0.55
1:A:10:PRO:HD3	1:A:27:ARG:HH12	1.72	0.54
1:B:17:VAL:O	1:B:20:GLU:HG2	2.07	0.54
1:A:211:TYR:CE1	1:A:235:ALA:HB2	2.43	0.54
1:B:114:THR:HG21	1:B:131:ARG:HH21	1.72	0.54
1:B:286:GLN:O	1:B:308:PRO:HA	2.07	0.54
1:A:80:ILE:N	1:A:80:ILE:CD1	2.70	0.53
1:B:113:VAL:HG12	1:B:115:ILE:HD12	1.90	0.53
1:A:8:GLU:O	1:A:27:ARG:NH1	2.41	0.53
1:A:182:ILE:N	1:A:182:ILE:CD1	2.71	0.53
1:B:251:LEU:HD23	1:B:252:VAL:N	2.23	0.53
1:B:260:GLY:HA3	1:B:279:GLY:H	1.73	0.53
1:A:180:ALA:HB1	1:A:182:ILE:HD11	1.90	0.52
1:B:289:ILE:HG13	1:B:311:ILE:HB	1.90	0.52
1:B:243:ASN:ND2	3:B:3138:HOH:O	2.43	0.52
1:A:10:PRO:HD3	1:A:27:ARG:NH1	2.24	0.52
1:A:102:ASP:O	1:A:103:ILE:HD12	2.09	0.52
1:A:274:HIS:HB2	1:A:291:GLY:O	2.09	0.52
1:B:92:ILE:CD1	1:B:92:ILE:N	2.66	0.52
1:A:63:ASP:OD2	1:A:65:ASN:HB2	2.09	0.52
1:B:81:THR:HB	1:B:82:GLN:NE2	2.25	0.52
1:B:251:LEU:C	1:B:251:LEU:CD2	2.77	0.52
1:A:211:TYR:CE1	1:A:235:ALA:CB	2.92	0.52
1:A:91:ARG:HD2	1:A:91:ARG:N	2.25	0.52
1:A:49:ALA:HB3	1:A:52:VAL:HG23	1.92	0.52
1:A:152:LEU:C	1:A:152:LEU:HD23	2.30	0.52
1:B:105:ASP:HB2	1:B:122:GLU:H	1.75	0.52
1:B:310:VAL:C	1:B:311:ILE:HD12	2.30	0.52
1:A:11:ARG:HG2	1:A:12:ALA:N	2.24	0.51
1:B:69:PHE:CD1	1:B:69:PHE:N	2.77	0.51
1:A:105:ASP:O	1:A:122:GLU:HA	2.10	0.51
1:B:190:ASP:OD2	1:B:211:TYR:HB2	2.10	0.51
1:A:73:GLU:CG	1:A:91:ARG:NH1	2.66	0.51
1:B:39:LYS:O	1:B:42:THR:HG23	2.11	0.51
1:B:85:THR:CG2	1:B:87:TYR:HE2	2.23	0.51
1:A:122:GLU:HB3	1:A:140:ALA:O	2.11	0.51
1:B:113:VAL:CG1	1:B:114:THR:N	2.72	0.51
1:B:152:LEU:HD11	1:B:165:ILE:HG22	1.92	0.51
1:B:170:GLN:O	1:B:187:GLU:HA	2.11	0.51
1:A:69:PHE:N	1:A:69:PHE:CD1	2.79	0.50
1:B:238:ALA:HB1	1:B:252:VAL:HG12	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:182:ILE:HD12	1:B:182:ILE:N	2.26	0.50
1:B:157:ARG:HH21	1:B:174:ASN:HD21	1.57	0.50
1:A:251:LEU:HD23	1:A:252:VAL:N	2.25	0.50
1:B:262:PRO:HD2	1:B:280:GLU:HG3	1.92	0.50
1:A:68:ALA:HA	1:A:86:LEU:O	2.12	0.50
1:A:34:ASP:OD1	1:A:40:ALA:N	2.32	0.49
1:A:314:GLU:HG2	1:B:97:TRP:CD2	2.47	0.49
1:B:262:PRO:CD	1:B:280:GLU:HG3	2.42	0.49
1:B:16:GLN:HA	1:B:20:GLU:O	2.13	0.49
1:A:100:ARG:O	1:A:100:ARG:HG2	2.12	0.49
1:B:239:LEU:HD23	1:B:240:ILE:N	2.27	0.49
1:A:90:VAL:HG21	1:A:103:ILE:HG22	1.95	0.48
1:A:113:VAL:CG1	1:A:114:THR:N	2.75	0.48
1:B:47:VAL:HA	1:B:68:ALA:O	2.12	0.48
1:B:165:ILE:N	1:B:165:ILE:CD1	2.74	0.48
1:A:103:ILE:HD12	1:A:103:ILE:N	2.27	0.48
1:A:166:VAL:HG12	1:A:167:HIS:N	2.28	0.48
1:B:298:PHE:O	1:B:299:ASP:C	2.51	0.48
1:B:311:ILE:N	1:B:311:ILE:CD1	2.76	0.48
1:B:96:VAL:HG21	1:B:109:ILE:HG22	1.95	0.48
1:B:274:HIS:CE1	3:B:3057:HOH:O	2.66	0.48
1:A:301:ASN:N	1:A:301:ASN:HD22	2.12	0.48
1:B:109:ILE:N	1:B:109:ILE:HD12	2.29	0.48
1:A:67:MSE:CE	1:A:85:THR:OG1	2.62	0.48
1:A:153:GLN:HB2	1:A:170:GLN:NE2	2.29	0.48
1:A:180:ALA:HB1	1:A:182:ILE:CD1	2.44	0.48
1:A:222:GLU:HG2	1:A:223:ASP:OD2	2.14	0.48
1:A:213:HIS:HE1	1:B:213:HIS:CE1	2.32	0.47
1:B:113:VAL:HG12	1:B:115:ILE:CD1	2.44	0.47
1:A:8:GLU:CA	3:A:3115:HOH:O	2.52	0.47
1:A:228:LEU:CD2	1:A:246:LEU:HD22	2.45	0.47
1:B:157:ARG:HG3	1:B:157:ARG:HH11	1.78	0.47
1:A:81:THR:O	1:A:82:GLN:HB2	2.13	0.47
1:A:103:ILE:N	1:A:103:ILE:CD1	2.77	0.47
1:B:164:ARG:C	1:B:165:ILE:HD12	2.35	0.47
1:A:85:THR:CG2	1:A:87:TYR:CD1	2.97	0.47
1:A:86:LEU:N	1:A:86:LEU:CD1	2.78	0.47
1:A:67:MSE:HE3	1:A:85:THR:OG1	2.15	0.47
1:A:113:VAL:HG12	1:A:114:THR:N	2.30	0.47
1:B:113:VAL:HG12	1:B:114:THR:N	2.28	0.47
1:A:126:ILE:HG23	1:A:154:ILE:HB	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:TYR:O	1:B:88:ASN:HB2	2.16	0.46
1:B:211:TYR:N	1:B:211:TYR:CD1	2.83	0.46
1:A:122:GLU:CB	1:A:140:ALA:O	2.64	0.45
1:A:306:ARG:HG2	1:A:306:ARG:HH11	1.81	0.45
1:A:274:HIS:CG	1:B:131:ARG:HD2	2.51	0.45
1:A:301:ASN:N	1:A:301:ASN:ND2	2.64	0.45
1:B:96:VAL:HG12	1:B:97:TRP:N	2.31	0.45
1:A:217:ILE:O	1:A:241:GLU:HA	2.16	0.45
1:A:301:ASN:CB	3:A:3126:HOH:O	2.63	0.45
1:A:15:TYR:CE2	1:A:17:VAL:CG2	3.00	0.45
1:A:8:GLU:HB3	3:A:3115:HOH:O	2.16	0.45
1:B:27:ARG:HH11	1:B:27:ARG:CG	2.25	0.45
1:A:11:ARG:HD2	1:A:13:PHE:CZ	2.52	0.45
1:B:26:LEU:HD22	1:B:47:VAL:C	2.38	0.45
1:B:204:ILE:CD1	1:B:204:ILE:N	2.80	0.44
1:B:248:HIS:HD2	3:B:3024:HOH:O	2.00	0.44
1:A:85:THR:CG2	1:A:87:TYR:HE1	2.29	0.44
1:B:194:ILE:HG23	1:B:202:VAL:HG11	2.00	0.44
1:A:176:THR:HG22	3:A:3066:HOH:O	2.17	0.44
1:B:169:VAL:HG13	1:B:186:ALA:O	2.16	0.44
1:A:239:LEU:HD23	1:A:240:ILE:N	2.33	0.43
1:B:39:LYS:NZ	1:B:39:LYS:CB	2.81	0.43
1:B:182:ILE:HG13	1:B:204:ILE:HB	2.00	0.43
1:A:33:THR:C	1:A:40:ALA:HB2	2.38	0.43
1:A:274:HIS:HE1	3:B:3033:HOH:O	2.01	0.43
1:B:81:THR:C	1:B:82:GLN:HE21	2.21	0.43
1:B:115:ILE:HD12	1:B:115:ILE:N	2.34	0.43
1:B:96:VAL:HG11	1:B:109:ILE:CG2	2.49	0.43
1:A:15:TYR:CE2	1:A:17:VAL:HG21	2.53	0.43
1:B:114:THR:HG21	1:B:131:ARG:NH2	2.34	0.43
1:A:15:TYR:CE1	1:A:22:LYS:HB3	2.54	0.43
1:A:243:ASN:ND2	3:A:3077:HOH:O	2.52	0.43
1:A:300:ASP:OD1	1:A:300:ASP:C	2.57	0.43
1:B:209:LYS:HD2	1:B:233:GLN:NE2	2.34	0.43
1:B:189:PHE:CD1	1:B:189:PHE:N	2.87	0.43
1:A:189:PHE:N	1:A:189:PHE:CD1	2.87	0.43
1:A:274:HIS:CE1	3:B:3033:HOH:O	2.72	0.43
1:B:74:ILE:HD12	1:B:74:ILE:N	2.33	0.43
1:A:121:ARG:O	1:A:122:GLU:CB	2.67	0.43
1:A:251:LEU:C	1:A:251:LEU:CD2	2.86	0.43
1:B:13:PHE:O	1:B:23:SER:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:96:VAL:CG2	1:B:109:ILE:HG22	2.48	0.43
1:B:268:ARG:HD2	3:B:3019:HOH:O	2.18	0.43
1:A:85:THR:HG22	1:A:87:TYR:HE1	1.78	0.42
1:B:96:VAL:CG1	1:B:97:TRP:N	2.82	0.42
1:B:157:ARG:HG3	1:B:157:ARG:NH1	2.34	0.42
1:A:122:GLU:CD	1:A:123:GLU:H	2.23	0.42
1:B:89:ASN:O	1:B:106:GLY:HA2	2.19	0.42
1:B:141:ILE:O	1:B:141:ILE:CG2	2.67	0.42
1:B:166:VAL:CG1	1:B:167:HIS:H	2.32	0.42
1:A:168:GLN:NE2	1:A:185:ARG:NH1	2.68	0.42
1:A:319:ARG:NH1	1:A:319:ARG:HG2	2.35	0.42
1:B:201:ASN:O	1:B:225:ILE:HG23	2.20	0.42
1:A:87:TYR:O	1:A:88:ASN:HB2	2.21	0.41
1:B:223:ASP:O	1:B:225:ILE:N	2.51	0.41
1:B:259:ARG:O	1:B:278:GLN:HA	2.20	0.41
1:A:15:TYR:CE1	1:A:22:LYS:CB	3.03	0.41
1:A:95:ASN:ND2	1:B:292:ARG:NH2	2.67	0.41
1:B:85:THR:CG2	1:B:87:TYR:CD2	3.03	0.41
1:B:169:VAL:HG22	1:B:182:ILE:HG22	2.01	0.41
1:B:169:VAL:CG1	1:B:170:GLN:N	2.83	0.41
1:B:285:ARG:HH12	1:B:308:PRO:CD	2.31	0.41
1:A:139:LEU:O	1:A:166:VAL:HA	2.20	0.41
1:A:316:ARG:NH1	3:A:3134:HOH:O	2.52	0.41
1:A:213:HIS:HB2	1:A:236:GLU:O	2.21	0.41
1:B:114:THR:C	1:B:115:ILE:HD12	2.41	0.41
1:A:104:SER:O	1:A:121:ARG:HA	2.21	0.41
1:B:85:THR:HG21	1:B:87:TYR:HE2	1.85	0.41
1:B:292:ARG:NH2	1:B:314:GLU:OE2	2.54	0.41
1:B:120:VAL:HG22	1:B:138:ILE:HD12	2.02	0.41
1:A:92:ILE:CD1	1:A:109:ILE:HD12	2.51	0.41
1:A:193:LEU:HD23	1:A:215:ARG:HG2	2.02	0.41
1:A:265:LEU:HD13	1:A:271:ILE:HD12	2.03	0.41
1:A:307:GLY:CA	1:A:308:PRO:C	2.88	0.41
1:B:184:HIS:HB3	1:B:185:ARG:H	1.62	0.41
1:B:217:ILE:O	1:B:241:GLU:HA	2.20	0.41
1:A:118:SER:HB3	1:A:136:SER:O	2.20	0.41
1:B:161:ASN:O	1:B:162:HIS:C	2.58	0.41
1:A:34:ASP:OD1	1:A:39:LYS:HA	2.22	0.40
1:A:89:ASN:O	1:A:106:GLY:HA2	2.21	0.40
1:A:180:ALA:CB	1:A:182:ILE:HD11	2.51	0.40
1:A:316:ARG:HD2	1:B:99:ASP:OD2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13:PHE:O	1:A:23:SER:HA	2.22	0.40
1:A:217:ILE:HD13	1:B:195:GLU:HG2	2.03	0.40
1:B:166:VAL:HG12	1:B:167:HIS:H	1.84	0.40
1:B:307:GLY:CA	1:B:308:PRO:C	2.89	0.40
1:B:59:CYS:HA	1:B:78:ALA:O	2.22	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	306/334 (92%)	273 (89%)	29 (10%)	4 (1%)	12	15
1	B	313/334 (94%)	283 (90%)	28 (9%)	2 (1%)	25	34
All	All	619/668 (93%)	556 (90%)	57 (9%)	6 (1%)	15	21

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	299	ASP
1	A	122	GLU
1	A	167	HIS
1	B	122	GLU
1	A	82	GLN
1	B	82	GLN

#### 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	255/275 (93%)	245 (96%)	10 (4%)	32	48
1	B	264/275 (96%)	250 (95%)	14 (5%)	22	34
All	All	519/550 (94%)	495 (95%)	24 (5%)	27	40

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	THR
1	A	56	GLN
1	A	91	ARG
1	A	133	LEU
1	A	182	ILE
1	A	198	LYS
1	A	243	ASN
1	A	285	ARG
1	A	300	ASP
1	A	319	ARG
1	B	16	GLN
1	B	27	ARG
1	B	39	LYS
1	B	82	GLN
1	B	91	ARG
1	B	110	SER
1	B	198	LYS
1	B	215	ARG
1	B	243	ASN
1	B	251	LEU
1	B	268	ARG
1	B	284	GLU
1	B	285	ARG
1	B	302	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	GLN
1	A	95	ASN
1	A	134	ASN

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Mol	Chain	Res	Type
1	A	135	GLN
1	A	150	GLN
1	A	168	GLN
1	A	170	GLN
1	A	237	HIS
1	A	243	ASN
1	A	248	HIS
1	A	255	HIS
1	A	274	HIS
1	A	301	ASN
1	B	56	GLN
1	B	82	GLN
1	B	134	ASN
1	B	135	GLN
1	B	168	GLN
1	B	170	GLN
1	B	237	HIS
1	B	243	ASN
1	B	248	HIS
1	B	255	HIS
1	B	274	HIS
1	B	278	GLN
1	B	301	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 monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	309/334 (92%)	0.25	17 (5%) 25 27	10, 21, 46, 60	0
1	B	318/334 (95%)	0.24	15 (4%) 31 34	12, 22, 38, 52	0
All	All	627/668 (93%)	0.24	32 (5%) 28 30	10, 22, 41, 60	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	299	ASP	7.3
1	A	300	ASP	5.0
1	A	21	LYS	4.2
1	B	141	ILE	4.1
1	A	50	ASP	4.1
1	B	299	ASP	3.9
1	B	16	GLN	3.8
1	A	122	GLU	3.8
1	B	181	PHE	3.7
1	B	203	TRP	3.5
1	A	319	ARG	3.5
1	A	16	GLN	3.4
1	B	56	GLN	3.4
1	A	149	ALA	3.1
1	B	298	PHE	3.1
1	B	61	ILE	2.8
1	A	298	PHE	2.8
1	B	50	ASP	2.7
1	A	260	GLY	2.7
1	B	260	GLY	2.7
1	A	20	GLU	2.6
1	B	21	LYS	2.5
1	A	11	ARG	2.4
1	B	47	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	11	ARG	2.3
1	B	122	GLU	2.3
1	A	301	ASN	2.3
1	A	19	GLY	2.2
1	B	19	GLY	2.1
1	A	12	ALA	2.1
1	A	39	LYS	2.1
1	A	18	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	ZN	A	335	1/1	0.99	0.04	34,34,34,34	0
2	ZN	A	336	1/1	0.99	0.09	44,44,44,44	0
2	ZN	B	335	1/1	0.99	0.09	41,41,41,41	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.