



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 07:44 AM UTC

PDB ID : 4E2C / pdb\_00004e2c  
Title : Crystal Structure of the periplasmic domain of the chimeric LPS O-antigen chain length regulator protein  
Authors : Kalynych, S.; Yao, D.; Magee, J.D.; Cygler, M.  
Deposited on : 2012-03-08  
Resolution : 2.80 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 : 4-5-2 with Phenix2.0  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

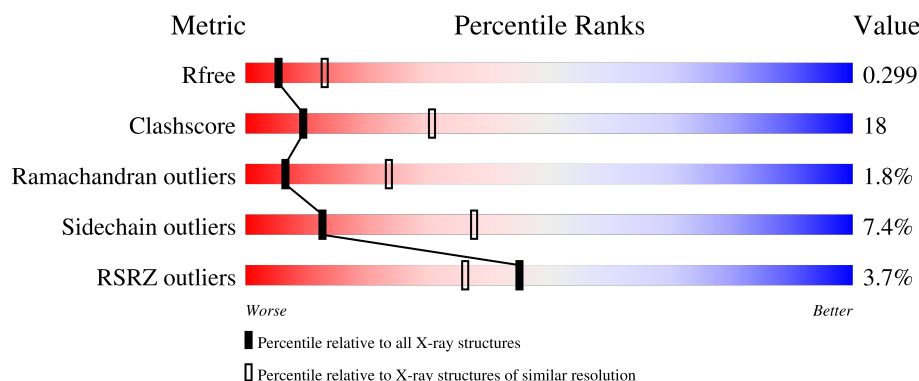
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
$R_{free}$	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	245	
1	B	245	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6823 atoms, of which 3238 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 chimeric WzzB Chain length determinant protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	233	Total	C	H	N	O	S	0	0	0
			3436	1113	1664	294	361	4			
1	B	232	Total	C	H	N	O	S	0	0	0
			3304	1088	1574	288	351	3			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	HIS	-	expression tag	UNP P37792
A	48	HIS	-	expression tag	UNP P37792
A	49	HIS	-	expression tag	UNP P37792
A	50	HIS	-	expression tag	UNP P37792
A	51	HIS	-	expression tag	UNP P37792
A	52	HIS	-	expression tag	UNP P37792
A	53	GLY	-	expression tag	UNP P37792
A	54	SER	-	expression tag	UNP P37792
B	47	HIS	-	expression tag	UNP P37792
B	48	HIS	-	expression tag	UNP P37792
B	49	HIS	-	expression tag	UNP P37792
B	50	HIS	-	expression tag	UNP P37792
B	51	HIS	-	expression tag	UNP P37792
B	52	HIS	-	expression tag	UNP P37792
B	53	GLY	-	expression tag	UNP P37792
B	54	SER	-	expression tag	UNP P37792

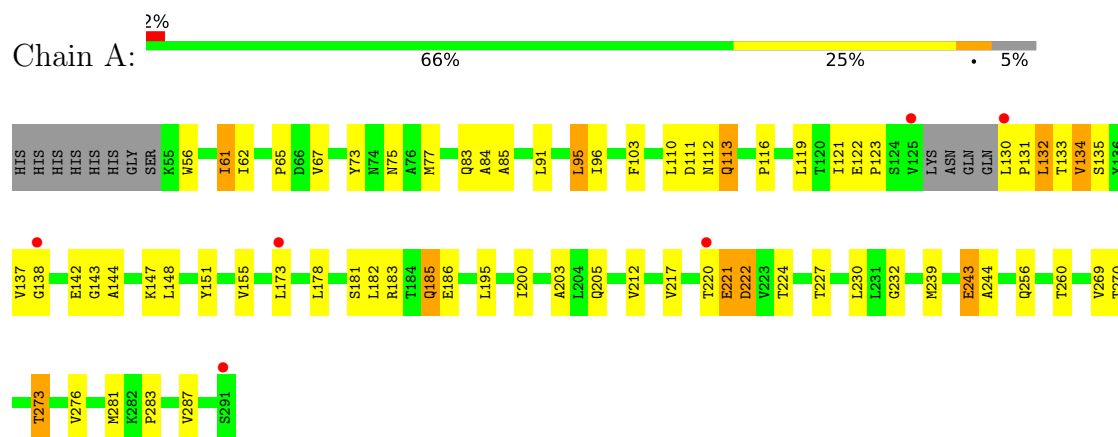
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	48	Total	O	0	0
			48	48		
2	B	35	Total	O	0	0
			35	35		

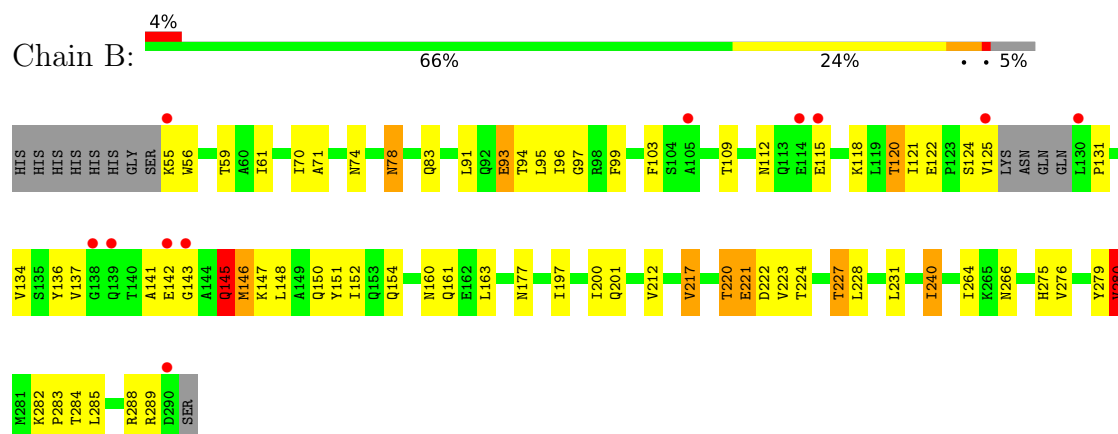
### 3 Residue-property plots [i](#)

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: chimeric WzzB Chain length determinant protein



- Molecule 1: chimeric WzzB Chain length determinant protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	116.95Å 116.95Å 219.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.50 – 2.80 42.50 – 2.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (42.50-2.80) 99.1 (42.50-2.80)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.71 (at 2.61Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.235 , 0.296 0.247 , 0.299	Depositor DCC
$R_{free}$ test set	1216 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	55.2	Xtriage
Anisotropy	0.237	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 56.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	6823	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.55% 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	0.60	1/1797 (0.1%)	0.90	4/2447 (0.2%)
1	B	0.59	1/1755 (0.1%)	0.88	5/2396 (0.2%)
All	All	0.60	2/3552 (0.1%)	0.89	9/4843 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	78	ASN	CG-ND2	-5.12	1.22	1.33
1	A	185	GLN	CD-NE2	-5.04	1.22	1.33

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	135	SER	N-CA-C	6.58	118.85	108.79
1	B	280	VAL	N-CA-C	-6.44	105.88	113.42
1	B	240	ILE	CB-CA-C	-5.59	104.69	112.02
1	A	130	LEU	CA-C-N	5.15	126.27	120.66
1	A	130	LEU	C-N-CA	5.15	126.27	120.66
1	B	212	VAL	CB-CA-C	-5.13	104.78	111.81
1	A	203	ALA	N-CA-C	-5.08	105.93	111.82
1	B	115	GLU	CA-C-N	5.04	124.80	119.76
1	B	115	GLU	C-N-CA	5.04	124.80	119.76

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	1772	1664	1664	60	2
1	B	1730	1574	1578	69	1
2	A	48	0	0	9	0
2	B	35	0	0	7	1
All	All	3585	3238	3242	121	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 18.

All (121) 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:142:GLU:OE2	1:B:285:LEU:CB	1.90	1.19
1:B:223:VAL:HG13	1:B:227:THR:HG22	1.39	1.03
1:A:121:ILE:HA	2:A:330:HOH:O	1.75	0.86
1:A:61:ILE:HG12	1:A:281:MET:HB3	1.59	0.83
1:A:270:THR:H	1:A:273:THR:CG2	1.95	0.80
2:A:343:HOH:O	1:B:227:THR:HG21	1.82	0.79
1:B:118:LYS:HB3	1:B:137:VAL:CG2	2.13	0.78
1:B:223:VAL:CG1	1:B:227:THR:HG22	2.13	0.76
1:B:118:LYS:HB3	1:B:137:VAL:HG22	1.67	0.75
1:A:205:GLN:O	2:A:308:HOH:O	2.08	0.70
1:A:220:THR:O	1:A:221:GLU:HB2	1.91	0.70
1:B:141:ALA:C	1:B:143:GLY:HA3	2.17	0.69
1:A:221:GLU:HA	1:A:221:GLU:OE1	1.91	0.69
1:A:270:THR:H	1:A:273:THR:HG23	1.59	0.68
1:B:103:PHE:HD2	1:B:121:ILE:CG1	2.06	0.67
1:B:223:VAL:HG13	1:B:227:THR:CG2	2.22	0.66
1:B:221:GLU:HG3	1:B:222:ASP:N	2.11	0.66
1:A:133:THR:O	2:A:330:HOH:O	2.13	0.65
1:B:142:GLU:N	1:B:143:GLY:HA3	2.10	0.65
1:B:124:SER:HB2	1:B:131:PRO:HG2	1.79	0.64
1:B:279:TYR:CD2	1:B:282:LYS:HG2	2.33	0.63
1:A:111:ASP:OD1	1:A:116:PRO:HB3	1.99	0.63
1:B:93:GLU:C	2:B:313:HOH:O	2.41	0.63
1:B:223:VAL:HG11	1:B:228:LEU:HA	1.80	0.63
1:B:220:THR:O	1:B:221:GLU:CB	2.46	0.62
1:A:200:ILE:HD13	1:A:239:MET:HE2	1.82	0.62
1:B:145:GLN:O	1:B:146:MET:C	2.44	0.60
1:A:269:VAL:HA	1:A:273:THR:HG21	1.83	0.60
1:B:197:ILE:O	1:B:201:GLN:HG3	2.02	0.58
1:A:56:TRP:O	1:A:138:GLY:N	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:ILE:CG1	1:A:281:MET:HB3	2.30	0.58
1:B:118:LYS:HB3	1:B:137:VAL:HG21	1.86	0.58
1:A:111:ASP:O	1:A:113:GLN:N	2.31	0.57
1:B:120:THR:HG23	1:B:122:GLU:OE1	2.03	0.57
1:A:103:PHE:CZ	1:A:119:LEU:HD11	2.40	0.57
1:B:220:THR:O	1:B:221:GLU:HB2	2.06	0.56
1:A:243:GLU:HG3	1:A:244:ALA:N	2.20	0.56
1:B:99:PHE:HB2	1:B:276:VAL:HG11	1.87	0.56
1:B:221:GLU:HG3	1:B:222:ASP:H	1.70	0.56
1:B:96:ILE:N	2:B:313:HOH:O	2.38	0.56
1:B:103:PHE:CD2	1:B:121:ILE:CG1	2.90	0.55
1:B:74:ASN:O	1:B:78:ASN:ND2	2.39	0.54
1:A:221:GLU:O	1:A:222:ASP:C	2.49	0.54
1:B:145:GLN:O	1:B:148:LEU:N	2.40	0.54
1:B:201:GLN:HG2	1:B:240:ILE:HD12	1.90	0.54
1:B:99:PHE:HB2	1:B:276:VAL:CG1	2.38	0.54
1:B:288:ARG:O	1:B:289:ARG:C	2.51	0.54
1:A:220:THR:O	1:A:221:GLU:CB	2.55	0.54
1:A:224:THR:OG1	1:A:227:THR:HG22	2.08	0.53
1:B:94:THR:C	2:B:313:HOH:O	2.52	0.53
1:B:151:TYR:O	1:B:152:ILE:C	2.51	0.52
1:A:138:GLY:HA3	1:A:144:ALA:HB2	1.90	0.52
1:B:61:ILE:O	1:B:280:VAL:HG13	2.09	0.52
1:A:222:ASP:HB2	1:B:224:THR:HG21	1.91	0.52
1:A:133:THR:HG22	2:A:346:HOH:O	2.09	0.52
1:A:131:PRO:C	1:A:132:LEU:HD13	2.35	0.52
1:A:270:THR:O	1:A:273:THR:HG23	2.10	0.52
1:B:96:ILE:HG22	2:B:313:HOH:O	2.09	0.51
1:A:122:GLU:HG2	2:A:338:HOH:O	2.09	0.51
1:A:110:LEU:HA	1:A:113:GLN:OE1	2.11	0.51
1:A:220:THR:CG2	1:B:217:VAL:HG21	2.41	0.51
1:A:137:VAL:HB	2:A:329:HOH:O	2.09	0.51
1:A:75:ASN:HD22	1:B:177:ASN:CG	2.19	0.50
1:B:95:LEU:N	2:B:313:HOH:O	2.44	0.50
1:B:148:LEU:HD23	1:B:283:PRO:HB3	1.93	0.50
1:B:282:LYS:HB3	1:B:283:PRO:HD2	1.94	0.50
1:B:222:ASP:O	1:B:222:ASP:OD1	2.30	0.49
1:A:220:THR:HG23	1:B:217:VAL:HG11	1.95	0.49
1:A:75:ASN:ND2	1:B:177:ASN:CG	2.71	0.49
1:B:125:VAL:HG12	1:B:125:VAL:O	2.13	0.48
1:B:160:ASN:O	1:B:161:GLN:C	2.56	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:GLN:O	1:B:154:GLN:HG3	2.13	0.48
1:A:103:PHE:CE2	1:A:134:VAL:HG22	2.49	0.48
1:A:148:LEU:HD23	1:A:283:PRO:HB3	1.95	0.48
1:B:95:LEU:HD13	1:B:275:HIS:O	2.14	0.47
1:A:121:ILE:HG13	2:A:330:HOH:O	2.14	0.47
1:A:182:LEU:O	1:A:186:GLU:HG3	2.14	0.47
1:B:55:LYS:O	1:B:289:ARG:HB2	2.14	0.47
1:B:142:GLU:CD	1:B:285:LEU:CB	2.81	0.47
1:A:111:ASP:C	1:A:113:GLN:N	2.72	0.47
1:A:73:TYR:CE1	1:A:77:MET:HE1	2.50	0.46
1:A:134:VAL:HA	2:A:330:HOH:O	2.14	0.46
1:B:223:VAL:CG1	1:B:228:LEU:HA	2.45	0.46
1:A:65:PRO:HG3	1:A:276:VAL:C	2.41	0.46
1:A:61:ILE:CD1	1:A:281:MET:HE2	2.45	0.46
1:A:200:ILE:CD1	1:A:239:MET:HE2	2.46	0.46
1:B:221:GLU:CG	1:B:222:ASP:N	2.77	0.46
1:A:181:SER:O	1:A:185:GLN:HG3	2.16	0.46
1:A:151:TYR:O	1:A:155:VAL:HG23	2.17	0.45
1:A:95:LEU:HD12	1:A:95:LEU:HA	1.82	0.45
1:B:70:ILE:O	1:B:71:ALA:C	2.59	0.45
1:A:212:VAL:HG12	1:A:232:GLY:HA2	1.98	0.45
1:B:56:TRP:CZ3	1:B:141:ALA:HB2	2.51	0.45
1:B:145:GLN:O	1:B:147:LYS:N	2.50	0.45
1:A:224:THR:H	1:A:227:THR:CG2	2.29	0.44
1:B:163:LEU:HD23	1:B:163:LEU:HA	1.87	0.44
1:A:143:GLY:O	1:A:147:LYS:HG3	2.18	0.44
1:A:217:VAL:HG13	1:A:230:LEU:HB3	2.00	0.44
1:B:96:ILE:HG23	1:B:97:GLY:N	2.33	0.44
1:A:123:PRO:HA	1:A:132:LEU:HD12	2.00	0.43
1:B:221:GLU:CG	1:B:222:ASP:H	2.31	0.43
1:A:83:GLN:H	1:A:83:GLN:CD	2.26	0.43
1:A:220:THR:HG22	1:B:217:VAL:HG21	1.99	0.43
1:A:287:VAL:HG23	2:B:304:HOH:O	2.18	0.43
1:B:200:ILE:HG22	1:B:240:ILE:HD11	1.99	0.43
1:B:266:ASN:HB3	2:B:307:HOH:O	2.17	0.43
1:B:121:ILE:HG23	1:B:134:VAL:HG12	2.01	0.43
1:A:103:PHE:CD2	1:A:121:ILE:HB	2.54	0.42
1:B:59:THR:O	1:B:284:THR:HG23	2.18	0.42
1:B:221:GLU:HA	1:B:221:GLU:OE1	2.19	0.42
1:A:61:ILE:HG12	1:A:281:MET:HE2	2.00	0.42
1:A:84:ALA:O	1:A:85:ALA:C	2.61	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:GLU:H	1:A:142:GLU:HG2	1.63	0.41
1:B:83:GLN:HA	1:B:83:GLN:OE1	2.20	0.41
1:A:178:LEU:HD11	1:A:260:THR:HG22	2.03	0.41
1:B:145:GLN:HB3	1:B:146:MET:H	1.64	0.41
1:A:221:GLU:O	1:A:222:ASP:O	2.39	0.41
1:A:220:THR:HG23	1:B:217:VAL:HG21	2.03	0.41
1:B:136:TYR:CE2	1:B:147:LYS:HB3	2.56	0.40
1:B:95:LEU:HD23	1:B:95:LEU:HA	1.81	0.40
1:A:220:THR:OG1	1:B:221:GLU:HB3	2.21	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)
1:A:221:GLU:O	1:B:220:THR:HG1[3_555]	1.58	0.02
1:A:227:THR:HG21	2:B:327:HOH:O[3_555]	1.58	0.02

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	229/245 (94%)	217 (95%)	9 (4%)	3 (1%)	9	31
1	B	228/245 (93%)	210 (92%)	13 (6%)	5 (2%)	5	19
All	All	457/490 (93%)	427 (93%)	22 (5%)	8 (2%)	6	23

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	221	GLU
1	B	145	GLN
1	B	146	MET

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Mol	Chain	Res	Type
1	B	221	GLU
1	B	220	THR
1	A	112	ASN
1	B	112	ASN
1	A	222	ASP

### 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	177/217 (82%)	162 (92%)	15 (8%)	10	31
1	B	163/217 (75%)	153 (94%)	10 (6%)	17	46
All	All	340/434 (78%)	315 (93%)	25 (7%)	13	37

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	ILE
1	A	62	ILE
1	A	67	VAL
1	A	91	LEU
1	A	95	LEU
1	A	96	ILE
1	A	113	GLN
1	A	132	LEU
1	A	134	VAL
1	A	173	LEU
1	A	183	ARG
1	A	195	LEU
1	A	243	GLU
1	A	256	GLN
1	A	273	THR
1	B	91	LEU
1	B	93	GLU
1	B	109	THR

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Mol	Chain	Res	Type
1	B	120	THR
1	B	145	GLN
1	B	217	VAL
1	B	227	THR
1	B	231	LEU
1	B	264	ILE
1	B	280	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	ASN
1	A	153	GLN
1	A	154	GLN
1	B	113	GLN
1	B	177	ASN
1	B	205	GLN
1	B	242	HIS
1	B	256	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 oligosaccharides 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, 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	233/245 (95%)	-0.12	6 (2%) 57 47	6, 25, 55, 114	0
1	B	232/245 (94%)	0.12	11 (4%) 36 29	4, 29, 76, 141	0
All	All	465/490 (94%)	0.00	17 (3%) 45 36	4, 26, 73, 141	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	291	SER	4.9
1	B	130	LEU	4.6
1	B	290	ASP	4.5
1	B	142	GLU	3.5
1	A	220	THR	3.3
1	B	105	ALA	3.1
1	B	114	GLU	3.1
1	B	139	GLN	2.9
1	A	130	LEU	2.8
1	A	125	VAL	2.6
1	B	115	GLU	2.6
1	A	138	GLY	2.5
1	B	138	GLY	2.5
1	B	55	LYS	2.4
1	B	125	VAL	2.2
1	B	143	GLY	2.2
1	A	173	LEU	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 oligosaccharides 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.