



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 03:16 am GMT

PDB ID : 2NLZ  
Title : Crystal structure of cephalosporin acylase from *Bacillus halodurans*  
Authors : Patskovsky, Y.; Ramagopal, U.; Sauder, J.M.; Dickey, M.; Adams, J.M.; Ozyurt, S.; Wasserman, S.R.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)  
Deposited on : 2006-10-20  
Resolution : 2.70 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
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

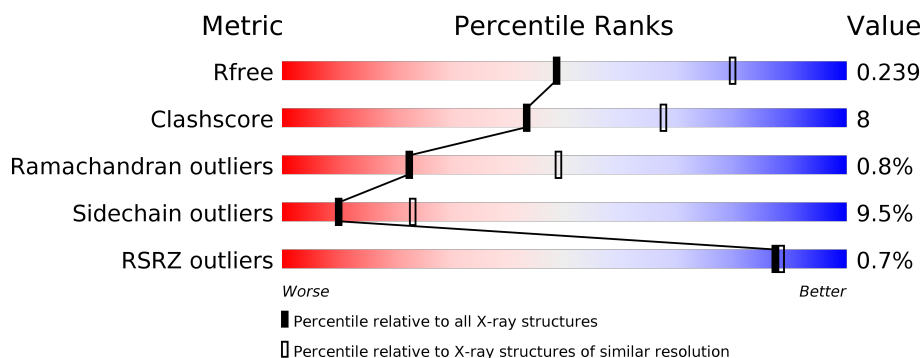
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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	2259 (2.70-2.70)
Clashscore	112137	2590 (2.70-2.70)
Ramachandran outliers	110173	2550 (2.70-2.70)
Sidechain outliers	110143	2550 (2.70-2.70)
RSRZ outliers	101464	2275 (2.70-2.70)

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	547	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>77%</span> <span>18%</span> <span>...</span> </div> </div>
1	B	547	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>77%</span> <span>18%</span> <span>...</span> </div> </div>
1	C	547	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>77%</span> <span>18%</span> <span>...</span> </div> </div>
1	D	547	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>78%</span> <span>18%</span> <span>...</span> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 16871 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 Cephalosporin acylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	536	Total	C	N	O	S	0	3	0
			4153	2645	713	778	17			
1	B	535	Total	C	N	O	S	0	1	0
			4137	2634	711	775	17			
1	C	535	Total	C	N	O	S	0	1	0
			4137	2634	711	775	17			
1	D	537	Total	C	N	O	S	0	2	0
			4163	2651	718	777	17			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	170	VAL	PHE	ENGINEERED	UNP Q9KEI5
A	540	GLU	-	CLONING ARTIFACT	UNP Q9KEI5
A	541	GLY	-	CLONING ARTIFACT	UNP Q9KEI5
A	542	HIS	-	EXPRESSION TAG	UNP Q9KEI5
A	543	HIS	-	EXPRESSION TAG	UNP Q9KEI5
A	544	HIS	-	EXPRESSION TAG	UNP Q9KEI5
A	545	HIS	-	EXPRESSION TAG	UNP Q9KEI5
A	546	HIS	-	EXPRESSION TAG	UNP Q9KEI5
A	547	HIS	-	EXPRESSION TAG	UNP Q9KEI5
B	170	VAL	PHE	ENGINEERED	UNP Q9KEI5
B	540	GLU	-	CLONING ARTIFACT	UNP Q9KEI5
B	541	GLY	-	CLONING ARTIFACT	UNP Q9KEI5
B	542	HIS	-	EXPRESSION TAG	UNP Q9KEI5
B	543	HIS	-	EXPRESSION TAG	UNP Q9KEI5
B	544	HIS	-	EXPRESSION TAG	UNP Q9KEI5
B	545	HIS	-	EXPRESSION TAG	UNP Q9KEI5
B	546	HIS	-	EXPRESSION TAG	UNP Q9KEI5
B	547	HIS	-	EXPRESSION TAG	UNP Q9KEI5
C	170	VAL	PHE	ENGINEERED	UNP Q9KEI5
C	540	GLU	-	CLONING ARTIFACT	UNP Q9KEI5
C	541	GLY	-	CLONING ARTIFACT	UNP Q9KEI5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	542	HIS	-	EXPRESSION TAG	UNP Q9KEI5
C	543	HIS	-	EXPRESSION TAG	UNP Q9KEI5
C	544	HIS	-	EXPRESSION TAG	UNP Q9KEI5
C	545	HIS	-	EXPRESSION TAG	UNP Q9KEI5
C	546	HIS	-	EXPRESSION TAG	UNP Q9KEI5
C	547	HIS	-	EXPRESSION TAG	UNP Q9KEI5
D	170	VAL	PHE	ENGINEERED	UNP Q9KEI5
D	540	GLU	-	CLONING ARTIFACT	UNP Q9KEI5
D	541	GLY	-	CLONING ARTIFACT	UNP Q9KEI5
D	542	HIS	-	EXPRESSION TAG	UNP Q9KEI5
D	543	HIS	-	EXPRESSION TAG	UNP Q9KEI5
D	544	HIS	-	EXPRESSION TAG	UNP Q9KEI5
D	545	HIS	-	EXPRESSION TAG	UNP Q9KEI5
D	546	HIS	-	EXPRESSION TAG	UNP Q9KEI5
D	547	HIS	-	EXPRESSION TAG	UNP Q9KEI5

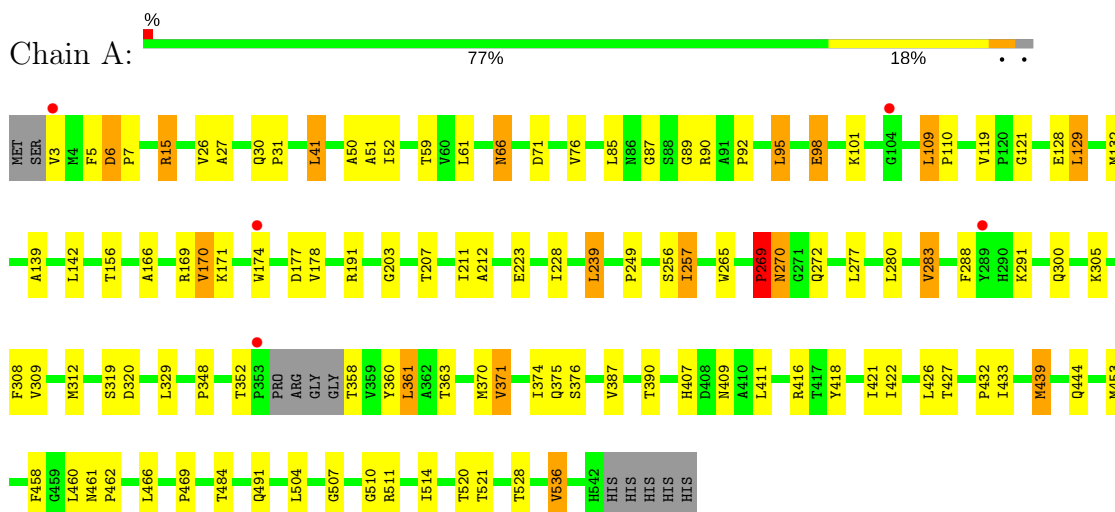
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	79	Total O 79 79	0	0
2	B	73	Total O 73 73	0	0
2	C	57	Total O 57 57	0	0
2	D	72	Total O 72 72	0	0

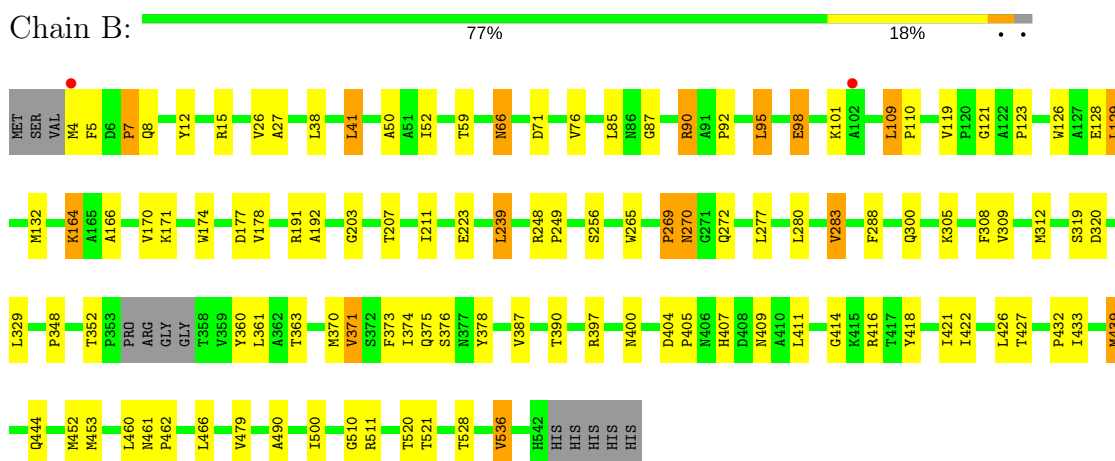
### 3 Residue-property plots [i](#)

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: Cephalosporin acylase

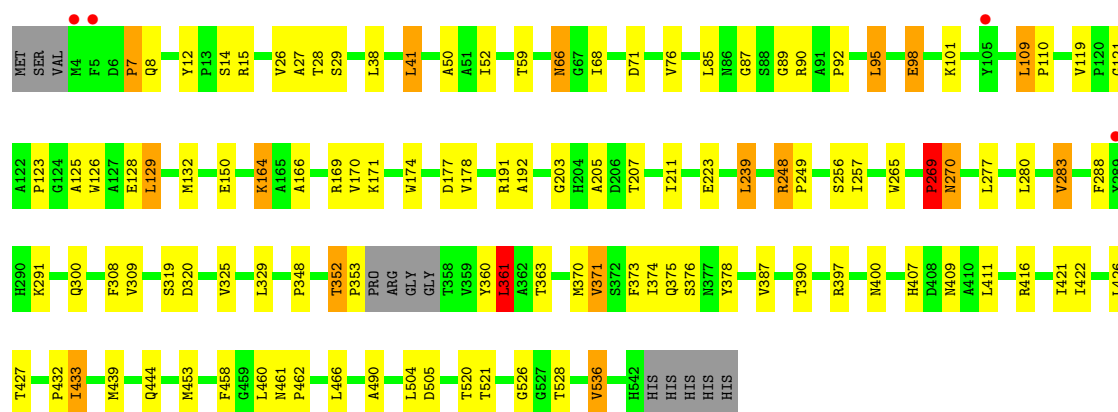


#### • Molecule 1: Cephalosporin acylase

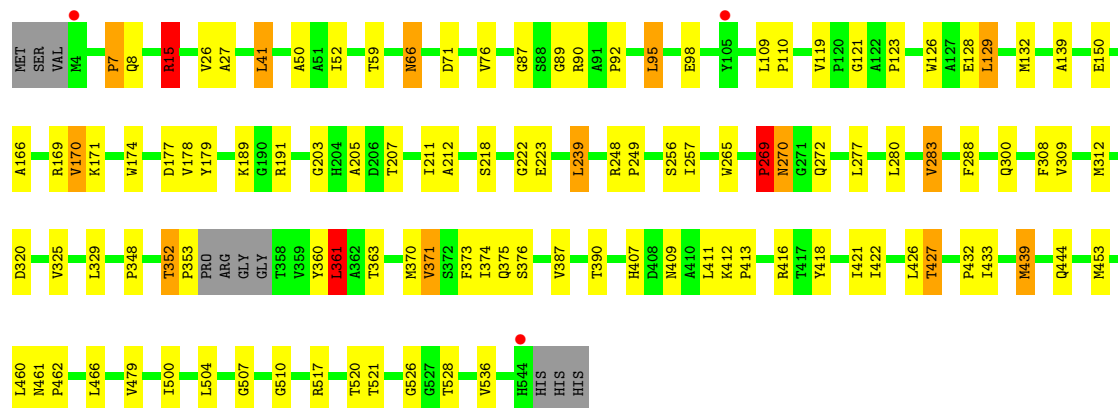
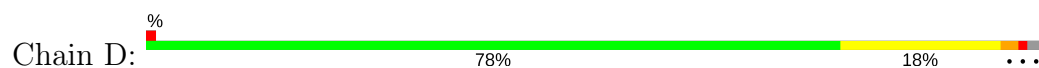


#### • Molecule 1: Cephalosporin acylase





● Molecule 1: Cephalosporin acylase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	105.75Å 105.75Å 385.06Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.70 19.98 – 2.70	Depositor EDS
% Data completeness (in resolution range)	95.2 (20.00-2.70) 94.9 (19.98-2.70)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.84 (at 2.71Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.210 , 0.245 0.209 , 0.239	Depositor DCC
$R_{free}$ test set	2011 reflections (3.13%)	DCC
Wilson B-factor (Å <sup>2</sup> )	37.5	Xtriage
Anisotropy	0.198	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 1.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.055 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	16871	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.40% 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

### 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	0.42	0/4277	0.61	2/5824 (0.0%)
1	B	0.40	0/4255	0.62	2/5794 (0.0%)
1	C	0.42	0/4255	0.61	1/5794 (0.0%)
1	D	0.41	0/4286	0.63	6/5835 (0.1%)
All	All	0.41	0/17073	0.62	11/23247 (0.0%)

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
1	B	0	1
1	C	0	1
1	D	0	1
All	All	0	4

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	361	LEU	CA-CB-CG	6.76	130.84	115.30
1	C	361	LEU	CA-CB-CG	6.52	130.30	115.30
1	D	517	ARG	NE-CZ-NH1	-6.39	117.10	120.30
1	A	361	LEU	CA-CB-CG	5.75	128.51	115.30
1	B	90	ARG	NE-CZ-NH2	-5.66	117.47	120.30
1	B	90	ARG	NE-CZ-NH1	5.66	123.13	120.30
1	D	15[A]	ARG	NE-CZ-NH2	5.58	123.09	120.30
1	D	15[B]	ARG	NE-CZ-NH2	5.58	123.09	120.30
1	D	15[A]	ARG	NE-CZ-NH1	-5.50	117.55	120.30
1	D	15[B]	ARG	NE-CZ-NH1	-5.50	117.55	120.30
1	A	142	LEU	CA-CB-CG	5.40	127.71	115.30



There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	269	PRO	Peptide
1	B	269	PRO	Peptide
1	C	269	PRO	Peptide
1	D	269	PRO	Peptide

## 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	4153	0	4035	68	0
1	B	4137	0	4013	62	0
1	C	4137	0	4013	65	0
1	D	4163	0	4040	63	0
2	A	79	0	0	3	0
2	B	73	0	0	2	0
2	C	57	0	0	2	0
2	D	72	0	0	1	0
All	All	16871	0	16101	251	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 8.

All (251) 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:376:SER:HB3	1:A:421:ILE:HD11	1.50	0.91
1:A:426:LEU:HD23	1:A:433:ILE:HD11	1.56	0.88
1:D:376:SER:HB3	1:D:421:ILE:HD11	1.59	0.84
1:B:376:SER:HB3	1:B:421:ILE:HD11	1.62	0.82
1:C:248:ARG:CG	1:C:248:ARG:HH11	1.93	0.82
1:C:248:ARG:NH1	1:C:248:ARG:HG2	1.95	0.82
1:C:166:ALA:O	1:C:170:VAL:HB	1.85	0.75
1:D:426:LEU:HD23	1:D:433:ILE:HD11	1.68	0.75
1:B:166:ALA:O	1:B:170:VAL:HB	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:THR:OG1	1:A:371:VAL:HG13	1.87	0.73
1:C:66:ASN:HD21	1:C:375:GLN:HG3	1.52	0.73
1:B:27:ALA:HB2	1:B:536:VAL:HG13	1.71	0.72
1:C:363:THR:OG1	1:C:371:VAL:HG13	1.89	0.72
1:C:407:HIS:HD2	1:C:409:ASN:H	1.37	0.71
1:D:66:ASN:HD21	1:D:375:GLN:HG3	1.55	0.71
1:D:166:ALA:O	1:D:170:VAL:HB	1.90	0.71
1:C:248:ARG:HH11	1:C:248:ARG:HG2	1.51	0.71
1:D:363:THR:OG1	1:D:371:VAL:HG13	1.92	0.70
1:B:461:ASN:HB2	1:B:462:PRO:HD2	1.74	0.70
1:A:66:ASN:HD21	1:A:375:GLN:HG3	1.57	0.69
1:B:407:HIS:HD2	1:B:409:ASN:H	1.38	0.69
1:A:166:ALA:O	1:A:170:VAL:HB	1.91	0.69
1:A:426:LEU:CD2	1:A:433:ILE:HD11	2.22	0.69
1:A:407:HIS:HD2	1:A:409:ASN:H	1.42	0.68
1:D:407:HIS:HD2	1:D:409:ASN:H	1.39	0.68
1:C:376:SER:HB3	1:C:421:ILE:HD11	1.76	0.68
1:B:426:LEU:HD23	1:B:433:ILE:HD11	1.76	0.67
1:B:363:THR:OG1	1:B:371:VAL:HG13	1.93	0.67
1:A:504:LEU:HD21	1:C:490:ALA:HB1	1.78	0.66
1:B:490:ALA:HB1	1:D:504:LEU:HD21	1.79	0.65
1:D:128:GLU:HG3	1:D:249:PRO:HG2	1.77	0.65
1:C:426:LEU:HD23	1:C:433:ILE:HD11	1.80	0.64
1:C:128:GLU:HG3	1:C:249:PRO:HG2	1.79	0.63
1:D:507:GLY:HA2	2:D:550:HOH:O	1.97	0.63
1:A:27:ALA:HB2	1:A:536:VAL:HG13	1.82	0.62
1:C:387:VAL:O	1:C:390:THR:HG22	1.98	0.62
1:C:12:TYR:O	1:D:15[B]:ARG:NH2	2.32	0.62
1:A:5:PHE:CZ	1:B:433:ILE:HD12	2.35	0.61
1:A:5:PHE:HZ	1:B:433:ILE:HD12	1.64	0.61
1:D:528:THR:HB	1:D:536:VAL:HG22	1.82	0.61
1:A:71:ASP:HB3	1:A:121:GLY:H	1.66	0.61
1:C:109:LEU:HB2	1:C:110:PRO:HD2	1.82	0.60
1:C:528:THR:HB	1:C:536:VAL:HG22	1.84	0.60
1:C:71:ASP:HB2	1:C:87:GLY:O	2.01	0.60
1:B:66:ASN:HD21	1:B:375:GLN:HG3	1.66	0.60
1:D:528:THR:HB	1:D:536:VAL:CG2	2.32	0.60
1:B:283:VAL:HG22	1:B:288:PHE:HZ	1.67	0.59
1:D:119:VAL:HG12	1:D:239:LEU:HD21	1.83	0.59
1:A:370:MET:CE	1:A:432:PRO:HB2	2.33	0.59
1:B:427:THR:HG22	2:B:605:HOH:O	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:300:GLN:HG2	1:D:453:MET:CE	2.32	0.59
1:A:109:LEU:HB2	1:A:110:PRO:HD2	1.84	0.59
1:C:309:VAL:HG21	1:C:348:PRO:HB2	1.83	0.59
1:A:528:THR:HB	1:A:536:VAL:HG22	1.85	0.59
1:B:422:ILE:H	1:B:444:GLN:HE22	1.51	0.58
1:C:71:ASP:HB3	1:C:121:GLY:H	1.68	0.58
1:C:300:GLN:HG2	1:C:453:MET:CE	2.33	0.58
1:C:461:ASN:HB2	1:C:462:PRO:HD2	1.85	0.58
1:A:156:THR:HG21	2:A:613:HOH:O	2.04	0.58
1:B:300:GLN:HG2	1:B:453:MET:HE1	1.86	0.58
1:A:5:PHE:C	1:A:6:ASP:OD1	2.42	0.57
1:B:119:VAL:HG12	1:B:239:LEU:HD21	1.85	0.57
1:A:376:SER:CB	1:A:421:ILE:HD11	2.30	0.57
1:A:128:GLU:HG3	1:A:249:PRO:HG2	1.87	0.57
1:D:300:GLN:HG2	1:D:453:MET:HE1	1.88	0.56
1:A:283:VAL:HG22	1:A:288:PHE:HZ	1.70	0.56
1:B:128:GLU:HG3	1:B:249:PRO:HG2	1.87	0.56
1:C:123:PRO:HA	1:C:126:TRP:CE3	2.40	0.56
1:D:374:ILE:HG23	1:D:421:ILE:HG23	1.88	0.56
1:B:376:SER:CB	1:B:421:ILE:HD11	2.34	0.56
1:D:387:VAL:O	1:D:390:THR:HG22	2.06	0.56
1:D:27:ALA:HB2	1:D:536:VAL:HG13	1.88	0.56
1:C:95:LEU:HD22	1:C:411:LEU:HD21	1.88	0.56
1:D:52:ILE:HD11	1:D:129:LEU:HB3	1.87	0.56
1:A:119:VAL:HG12	1:A:239:LEU:HD21	1.88	0.56
1:D:283:VAL:HG22	1:D:288:PHE:HZ	1.71	0.56
1:B:71:ASP:HB2	1:B:87:GLY:O	2.06	0.55
1:C:52:ILE:HD11	1:C:129:LEU:HB3	1.87	0.55
1:D:370:MET:CE	1:D:432:PRO:HB2	2.37	0.55
1:D:71:ASP:HB3	1:D:121:GLY:H	1.70	0.55
1:B:300:GLN:HG2	1:B:453:MET:CE	2.35	0.55
1:B:90:ARG:HD3	1:B:414:GLY:O	2.07	0.54
1:C:374:ILE:HG23	1:C:421:ILE:HG23	1.89	0.54
1:D:109:LEU:HB2	1:D:110:PRO:HD2	1.89	0.54
1:D:26:VAL:HG22	1:D:363:THR:HG22	1.90	0.54
1:A:26:VAL:HG22	1:A:363:THR:HG22	1.89	0.54
1:B:265:TRP:HE1	1:B:427:THR:HB	1.72	0.54
1:C:397:ARG:HD3	1:C:400:ASN:HD22	1.73	0.54
1:B:109:LEU:HB2	1:B:110:PRO:HD2	1.90	0.54
1:D:374:ILE:HG23	1:D:421:ILE:CG2	2.38	0.53
1:B:370:MET:CE	1:B:432:PRO:HB2	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:387:VAL:O	1:A:390:THR:HG22	2.08	0.53
1:A:439:MET:HE3	1:A:510:GLY:HA2	1.91	0.53
1:D:426:LEU:CD2	1:D:433:ILE:HD11	2.36	0.53
1:A:461:ASN:HB2	1:A:462:PRO:HD2	1.91	0.53
1:A:5:PHE:O	1:A:6:ASP:OD1	2.27	0.53
1:B:95:LEU:HD22	1:B:411:LEU:HD21	1.90	0.53
1:B:92:PRO:HG3	1:B:239:LEU:HD23	1.90	0.53
1:C:26:VAL:HG22	1:C:363:THR:HG22	1.91	0.53
1:D:376:SER:CB	1:D:421:ILE:HD11	2.34	0.53
1:C:128:GLU:HG2	2:C:579:HOH:O	2.09	0.52
1:B:123:PRO:HA	1:B:126:TRP:CE3	2.44	0.52
1:B:528:THR:HB	1:B:536:VAL:HG22	1.92	0.52
1:B:71:ASP:HB3	1:B:121:GLY:H	1.74	0.52
1:C:374:ILE:HG23	1:C:421:ILE:CG2	2.40	0.52
1:B:309:VAL:HG21	1:B:348:PRO:HB2	1.91	0.51
1:D:170:VAL:CG1	1:D:174:TRP:CE3	2.94	0.51
1:B:92:PRO:HD2	1:B:95:LEU:HD12	1.92	0.51
1:C:361:LEU:HD22	1:C:373:PHE:HB3	1.91	0.51
1:A:439:MET:HB3	1:A:511:ARG:H	1.76	0.51
1:B:387:VAL:O	1:B:390:THR:HG22	2.10	0.51
1:B:511:ARG:NH2	2:B:549:HOH:O	2.43	0.51
1:C:269:PRO:HG2	1:C:325:VAL:HG22	1.93	0.51
1:B:203:GLY:O	1:B:207:THR:HG22	2.11	0.50
1:C:505:ASP:HB3	2:C:571:HOH:O	2.11	0.50
1:B:361:LEU:HD22	1:B:373:PHE:HB3	1.92	0.50
1:A:309:VAL:HG21	1:A:348:PRO:HB2	1.94	0.50
1:B:52:ILE:HD11	1:B:129:LEU:HB3	1.94	0.49
1:C:265:TRP:HE1	1:C:427:THR:HB	1.76	0.49
1:C:370:MET:CE	1:C:432:PRO:HB2	2.42	0.49
1:A:89:GLY:HA3	1:A:119:VAL:HG23	1.95	0.49
1:C:89:GLY:HA3	1:C:119:VAL:HG23	1.95	0.49
1:B:41:LEU:HD13	1:B:50:ALA:HB2	1.95	0.49
1:D:439:MET:HE3	1:D:510:GLY:HA2	1.95	0.49
1:D:170:VAL:HG11	1:D:174:TRP:CE3	2.47	0.49
1:D:123:PRO:HA	1:D:126:TRP:CE3	2.48	0.49
1:D:374:ILE:CG2	1:D:421:ILE:HG23	2.42	0.49
1:D:461:ASN:HB2	1:D:462:PRO:HD2	1.96	0.48
1:A:300:GLN:HG2	1:A:453:MET:CE	2.43	0.48
1:C:119:VAL:HG12	1:C:239:LEU:HD21	1.94	0.48
1:D:256:SER:HB3	1:D:265:TRP:CE3	2.48	0.48
1:B:256:SER:HB3	1:B:265:TRP:CE3	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:305:LYS:O	1:A:309:VAL:HG23	2.14	0.48
1:A:507:GLY:HA2	2:A:604:HOH:O	2.14	0.48
1:A:95:LEU:HD22	1:A:411:LEU:HD21	1.96	0.48
1:B:426:LEU:CD2	1:B:433:ILE:HD11	2.43	0.48
1:C:27:ALA:HB2	1:C:536:VAL:HG13	1.95	0.48
1:D:265:TRP:HE1	1:D:427:THR:HB	1.79	0.47
1:D:41:LEU:HD13	1:D:50:ALA:HB2	1.94	0.47
1:A:257:ILE:HD13	1:A:257:ILE:HA	1.80	0.47
1:D:272:GLN:NE2	1:D:418:TYR:OH	2.48	0.47
1:A:203:GLY:O	1:A:207:THR:HG22	2.14	0.47
1:C:422:ILE:H	1:C:444:GLN:HE22	1.63	0.47
1:C:92:PRO:HD2	1:C:95:LEU:HD12	1.96	0.47
1:B:170:VAL:HG13	1:B:174:TRP:CD2	2.50	0.47
1:D:352:THR:HB	1:D:353:PRO:HD2	1.96	0.47
1:C:376:SER:CB	1:C:421:ILE:HD11	2.43	0.46
1:A:109:LEU:HD23	2:A:625:HOH:O	2.14	0.46
1:A:265:TRP:HE1	1:A:427:THR:HB	1.81	0.46
1:D:170:VAL:HG13	1:D:174:TRP:CD2	2.51	0.46
1:A:71:ASP:HB2	1:A:87:GLY:O	2.16	0.46
1:D:361:LEU:HD22	1:D:373:PHE:HB3	1.97	0.46
1:A:41:LEU:HD13	1:A:50:ALA:HB2	1.98	0.46
1:A:85:LEU:HD11	1:A:249:PRO:HB2	1.98	0.46
1:A:374:ILE:HG23	1:A:421:ILE:HG23	1.98	0.46
1:C:85:LEU:HD11	1:C:249:PRO:HB2	1.97	0.46
1:A:52:ILE:HD11	1:A:129:LEU:HB3	1.97	0.46
1:B:374:ILE:HG23	1:B:421:ILE:CG2	2.46	0.46
1:D:95:LEU:HD22	1:D:411:LEU:HD21	1.98	0.45
1:A:291:LYS:HB2	1:A:458:PHE:CG	2.52	0.45
1:D:150:GLU:HG3	1:D:205:ALA:CB	2.47	0.45
1:A:228:ILE:HG12	1:A:387:VAL:HG21	1.99	0.45
1:B:85:LEU:HD11	1:B:249:PRO:HB2	1.98	0.45
1:A:51:ALA:HA	1:A:361:LEU:HD21	1.99	0.45
1:D:139:ALA:HA	1:D:212:ALA:HB1	1.99	0.45
1:C:426:LEU:CD2	1:C:433:ILE:HD11	2.46	0.45
1:D:309:VAL:HG21	1:D:348:PRO:HB2	1.98	0.45
1:A:374:ILE:HG23	1:A:421:ILE:CG2	2.48	0.44
1:C:256:SER:HB3	1:C:265:TRP:CE3	2.53	0.44
1:D:526:GLY:HA3	1:D:536:VAL:HG11	1.98	0.44
1:A:272:GLN:NE2	1:A:418:TYR:OH	2.51	0.44
1:A:376:SER:HB3	1:A:421:ILE:CD1	2.36	0.44
1:B:170:VAL:CG1	1:B:174:TRP:CE3	3.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:421:ILE:HG21	1:C:421:ILE:HD13	1.66	0.44
1:B:374:ILE:HG23	1:B:421:ILE:HG23	1.99	0.44
1:A:66:ASN:HD22	1:A:66:ASN:HA	1.63	0.44
1:C:170:VAL:HG13	1:C:174:TRP:CD2	2.53	0.44
1:A:170:VAL:CG1	1:A:174:TRP:CE3	3.01	0.44
1:C:283:VAL:HG22	1:C:288:PHE:HZ	1.83	0.44
1:C:125:ALA:O	1:C:129:LEU:HB2	2.18	0.44
1:B:26:VAL:HG22	1:B:363:THR:HG22	2.00	0.43
1:B:397:ARG:HD3	1:B:400:ASN:HD22	1.84	0.43
1:D:71:ASP:HB2	1:D:87:GLY:O	2.17	0.43
1:B:404:ASP:HA	1:B:405:PRO:HD3	1.89	0.43
1:A:491[B]:GLN:HG2	1:C:504:LEU:HG	2.00	0.43
1:A:370:MET:HE3	1:A:432:PRO:HB2	2.00	0.43
1:B:98:GLU:HA	1:B:101:LYS:HB3	2.00	0.43
1:C:98:GLU:HA	1:C:101:LYS:HB3	2.01	0.43
1:D:170:VAL:HG13	1:D:174:TRP:CE3	2.54	0.43
1:D:363:THR:HG1	1:D:371:VAL:HG13	1.80	0.43
1:D:421:ILE:HD13	1:D:421:ILE:HG21	1.74	0.43
1:A:139:ALA:HA	1:A:212:ALA:HB1	2.00	0.43
1:C:291:LYS:HB2	1:C:458:PHE:CG	2.54	0.43
1:D:479:VAL:HB	1:D:500:ILE:HD13	2.01	0.43
1:A:358:THR:HG23	1:A:376:SER:HB2	2.01	0.43
1:B:164:LYS:HG2	1:B:192:ALA:HB3	2.00	0.43
1:C:352:THR:HB	1:C:353:PRO:HD2	2.01	0.43
1:D:412:LYS:HG3	1:D:413:PRO:HD2	2.01	0.43
1:C:374:ILE:CG2	1:C:421:ILE:HG23	2.48	0.42
1:A:170:VAL:HG13	1:A:174:TRP:CD2	2.55	0.42
1:A:528:THR:HB	1:A:536:VAL:CG2	2.47	0.42
1:B:452:MET:HA	1:B:452:MET:HE2	2.00	0.42
1:D:92:PRO:HD2	1:D:95:LEU:HD12	2.01	0.42
1:B:256:SER:HB3	1:B:265:TRP:CD2	2.55	0.42
1:A:98:GLU:HA	1:A:101:LYS:HB3	2.01	0.42
1:B:479:VAL:HB	1:B:500:ILE:HD13	2.01	0.42
1:C:41:LEU:HD13	1:C:50:ALA:HB2	2.01	0.42
1:D:92:PRO:HG3	1:D:239:LEU:HD23	2.00	0.42
1:B:360:TYR:CG	1:B:361:LEU:N	2.87	0.42
1:B:7:PRO:HB2	1:B:8:GLN:H	1.52	0.42
1:D:352:THR:HB	1:D:353:PRO:CD	2.49	0.42
1:A:170:VAL:HG11	1:A:174:TRP:CE3	2.54	0.42
1:A:256:SER:HB3	1:A:265:TRP:CE3	2.54	0.42
1:A:469:PRO:HB3	1:A:484:THR:HG23	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:421:ILE:HG21	1:B:421:ILE:HD13	1.81	0.42
1:C:28:THR:OG1	1:C:29:SER:N	2.52	0.42
1:C:170:VAL:CG1	1:C:174:TRP:CE3	3.03	0.42
1:C:164:LYS:HG2	1:C:192:ALA:HB3	2.01	0.41
1:C:207:THR:HG21	1:C:390:THR:O	2.20	0.41
1:A:360:TYR:CG	1:A:361:LEU:N	2.88	0.41
1:B:422:ILE:H	1:B:444:GLN:NE2	2.16	0.41
1:B:422:ILE:HG13	1:B:422:ILE:O	2.21	0.41
1:C:170:VAL:HG13	1:C:174:TRP:CE3	2.55	0.41
1:C:528:THR:HB	1:C:536:VAL:CG2	2.50	0.41
1:B:272:GLN:NE2	1:B:418:TYR:OH	2.54	0.41
1:C:300:GLN:HG2	1:C:453:MET:HE1	2.01	0.41
1:D:218:SER:HA	1:D:222:GLY:HA3	2.02	0.41
1:B:4:MET:HB3	1:B:5:PHE:H	1.70	0.41
1:C:66:ASN:HD22	1:C:66:ASN:HA	1.64	0.41
1:C:150:GLU:HG3	1:C:205:ALA:CB	2.50	0.41
1:C:360:TYR:CG	1:C:361:LEU:N	2.88	0.41
1:D:360:TYR:CG	1:D:361:LEU:N	2.89	0.41
1:D:7:PRO:HB2	1:D:8:GLN:H	1.56	0.41
1:A:15[B]:ARG:NH2	1:B:12:TYR:O	2.54	0.41
1:D:109:LEU:HD23	1:D:109:LEU:H	1.86	0.41
1:D:269:PRO:HG2	1:D:325:VAL:HG22	2.01	0.41
1:A:30:GLN:HG2	1:A:61:LEU:HB3	2.02	0.41
1:D:422:ILE:H	1:D:444:GLN:HE22	1.69	0.41
1:A:462:PRO:HB3	1:A:514:ILE:CG2	2.51	0.40
1:C:7:PRO:HB2	1:C:8:GLN:H	1.58	0.40
1:A:30:GLN:HA	1:A:31:PRO:HD3	1.95	0.40
1:A:92:PRO:HG3	1:A:239:LEU:HD23	2.03	0.40
1:A:92:PRO:HD2	1:A:95:LEU:HD12	2.03	0.40
1:C:203:GLY:O	1:C:207:THR:HG22	2.21	0.40
1:C:526:GLY:HA3	1:C:536:VAL:HG11	2.02	0.40
1:D:174:TRP:HB3	1:D:179:TYR:CG	2.55	0.40
1:D:203:GLY:O	1:D:207:THR:HG22	2.20	0.40
1:B:305:LYS:O	1:B:309:VAL:HG23	2.21	0.40
1:D:89:GLY:HA3	1:D:119:VAL:HG23	2.02	0.40
1:A:422:ILE:H	1:A:444:GLN:HE22	1.69	0.40
1:B:439:MET:HE3	1:B:510:GLY:HA2	2.04	0.40

There are no symmetry-related clashes.

## 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	535/547 (98%)	509 (95%)	22 (4%)	4 (1%)	25	53
1	B	532/547 (97%)	511 (96%)	16 (3%)	5 (1%)	20	46
1	C	532/547 (97%)	506 (95%)	21 (4%)	5 (1%)	20	46
1	D	535/547 (98%)	509 (95%)	22 (4%)	4 (1%)	25	53
All	All	2134/2188 (98%)	2035 (95%)	81 (4%)	18 (1%)	22	49

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	439	MET
1	C	439	MET
1	D	439	MET
1	A	7	PRO
1	B	7	PRO
1	C	269	PRO
1	C	270	ASN
1	D	269	PRO
1	A	269	PRO
1	A	270	ASN
1	B	269	PRO
1	B	270	ASN
1	C	7	PRO
1	D	7	PRO
1	D	270	ASN
1	B	378	TYR
1	B	439	MET
1	C	378	TYR



### 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	428/434 (99%)	385 (90%)	43 (10%)	9	21
1	B	425/434 (98%)	387 (91%)	38 (9%)	11	26
1	C	425/434 (98%)	381 (90%)	44 (10%)	8	18
1	D	428/434 (99%)	386 (90%)	42 (10%)	9	21
All	All	1706/1736 (98%)	1539 (90%)	167 (10%)	10	21

All (167) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	VAL
1	A	6	ASP
1	A	15[A]	ARG
1	A	15[B]	ARG
1	A	41	LEU
1	A	59	THR
1	A	66	ASN
1	A	76	VAL
1	A	90	ARG
1	A	95	LEU
1	A	98	GLU
1	A	109	LEU
1	A	129	LEU
1	A	132	MET
1	A	169	ARG
1	A	170	VAL
1	A	171	LYS
1	A	177	ASP
1	A	178	VAL
1	A	191	ARG
1	A	211	ILE
1	A	223	GLU
1	A	239	LEU
1	A	257	ILE

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Mol	Chain	Res	Type
1	A	269	PRO
1	A	270	ASN
1	A	277	LEU
1	A	280	LEU
1	A	283	VAL
1	A	308	PHE
1	A	312	MET
1	A	319[A]	SER
1	A	319[B]	SER
1	A	320	ASP
1	A	329	LEU
1	A	352	THR
1	A	371	VAL
1	A	416	ARG
1	A	460	LEU
1	A	466	LEU
1	A	520	THR
1	A	521	THR
1	A	536	VAL
1	B	15[A]	ARG
1	B	15[B]	ARG
1	B	38	LEU
1	B	41	LEU
1	B	59	THR
1	B	66	ASN
1	B	76	VAL
1	B	95	LEU
1	B	98	GLU
1	B	109	LEU
1	B	129	LEU
1	B	132	MET
1	B	164	LYS
1	B	171	LYS
1	B	177	ASP
1	B	178	VAL
1	B	191	ARG
1	B	211	ILE
1	B	223	GLU
1	B	239	LEU
1	B	248	ARG
1	B	270	ASN
1	B	277	LEU

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Mol	Chain	Res	Type
1	B	280	LEU
1	B	283	VAL
1	B	308	PHE
1	B	312	MET
1	B	319	SER
1	B	320	ASP
1	B	329	LEU
1	B	352	THR
1	B	371	VAL
1	B	416	ARG
1	B	460	LEU
1	B	466	LEU
1	B	520	THR
1	B	521	THR
1	B	536	VAL
1	C	14	SER
1	C	15[A]	ARG
1	C	15[B]	ARG
1	C	38	LEU
1	C	41	LEU
1	C	59	THR
1	C	66	ASN
1	C	68	ILE
1	C	76	VAL
1	C	90	ARG
1	C	95	LEU
1	C	98	GLU
1	C	109	LEU
1	C	129	LEU
1	C	132	MET
1	C	164	LYS
1	C	169	ARG
1	C	171	LYS
1	C	177	ASP
1	C	178	VAL
1	C	191	ARG
1	C	211	ILE
1	C	223	GLU
1	C	239	LEU
1	C	248	ARG
1	C	257	ILE
1	C	270	ASN

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Mol	Chain	Res	Type
1	C	277	LEU
1	C	280	LEU
1	C	283	VAL
1	C	308	PHE
1	C	319	SER
1	C	320	ASP
1	C	329	LEU
1	C	352	THR
1	C	361	LEU
1	C	371	VAL
1	C	416	ARG
1	C	433	ILE
1	C	460	LEU
1	C	466	LEU
1	C	520	THR
1	C	521	THR
1	C	536	VAL
1	D	15[A]	ARG
1	D	15[B]	ARG
1	D	41	LEU
1	D	59	THR
1	D	66	ASN
1	D	76	VAL
1	D	90	ARG
1	D	95	LEU
1	D	98	GLU
1	D	129	LEU
1	D	132	MET
1	D	169	ARG
1	D	170	VAL
1	D	171[A]	LYS
1	D	171[B]	LYS
1	D	177	ASP
1	D	178	VAL
1	D	189	LYS
1	D	191	ARG
1	D	211	ILE
1	D	223	GLU
1	D	239	LEU
1	D	248	ARG
1	D	257	ILE
1	D	269	PRO

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Mol	Chain	Res	Type
1	D	270	ASN
1	D	277	LEU
1	D	280	LEU
1	D	283	VAL
1	D	308	PHE
1	D	312	MET
1	D	320	ASP
1	D	329	LEU
1	D	352	THR
1	D	361	LEU
1	D	371	VAL
1	D	416	ARG
1	D	427	THR
1	D	460	LEU
1	D	466	LEU
1	D	520	THR
1	D	521	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (46) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	66	ASN
1	A	272	GLN
1	A	290	HIS
1	A	400	ASN
1	A	406	ASN
1	A	407	HIS
1	A	444	GLN
1	A	478	GLN
1	A	501	GLN
1	A	513	GLN
1	B	23	ASN
1	B	35	GLN
1	B	66	ASN
1	B	272	GLN
1	B	400	ASN
1	B	406	ASN
1	B	407	HIS
1	B	444	GLN
1	B	472	GLN
1	B	501	GLN

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Mol	Chain	Res	Type
1	B	513	GLN
1	C	23	ASN
1	C	35	GLN
1	C	66	ASN
1	C	272	GLN
1	C	400	ASN
1	C	406	ASN
1	C	407	HIS
1	C	444	GLN
1	C	478	GLN
1	C	501	GLN
1	C	513	GLN
1	D	23	ASN
1	D	30	GLN
1	D	66	ASN
1	D	272	GLN
1	D	290	HIS
1	D	375	GLN
1	D	400	ASN
1	D	407	HIS
1	D	444	GLN
1	D	478	GLN
1	D	501	GLN
1	D	513	GLN
1	D	542	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

There are no ligands in this entry.

## 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 [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	536/547 (97%)	-0.51	5 (0%) 84 85	11, 39, 93, 162	0
1	B	535/547 (97%)	-0.51	2 (0%) 92 93	11, 38, 93, 162	0
1	C	535/547 (97%)	-0.51	4 (0%) 87 88	11, 39, 93, 162	0
1	D	537/547 (98%)	-0.52	3 (0%) 89 90	12, 39, 94, 161	0
All	All	2143/2188 (97%)	-0.51	14 (0%) 87 88	11, 39, 93, 162	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	4	MET	4.2
1	A	289	TYR	4.0
1	C	289	TYR	2.9
1	D	4	MET	2.9
1	D	105	TYR	2.9
1	B	102	ALA	2.7
1	A	174	TRP	2.7
1	A	104	GLY	2.6
1	B	4	MET	2.6
1	A	353	PRO	2.4
1	D	544	HIS	2.3
1	C	105	TYR	2.2
1	C	5	PHE	2.2
1	A	3	VAL	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.