



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:30 PM GMT

PDB ID : 4CDI  
Title : Crystal structure of AcrB-AcrZ complex  
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Deposited on : 2013-10-31  
Resolution : 3.70 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  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

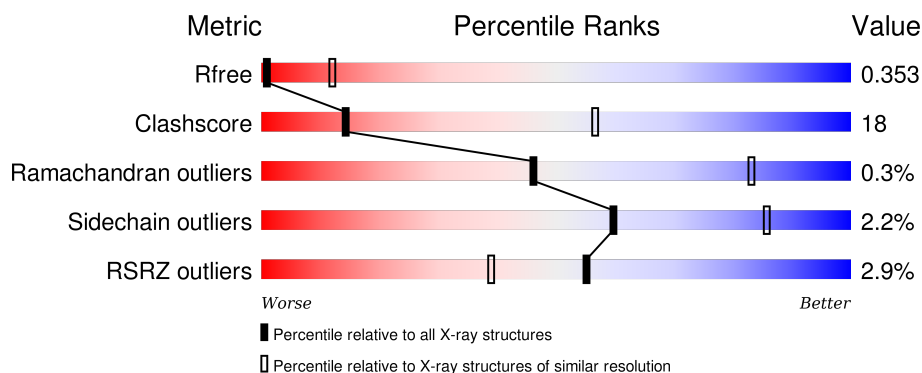
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.70 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))
$R_{free}$	91344	1101 (3.90-3.50)
Clashscore	102246	1224 (3.90-3.50)
Ramachandran outliers	100387	1172 (3.90-3.50)
Sidechain outliers	100360	1170 (3.90-3.50)
RSRZ outliers	91569	1108 (3.90-3.50)

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	1049	
2	C	49	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8114 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 ACRIFLAVINE RESISTANCE PROTEIN B.

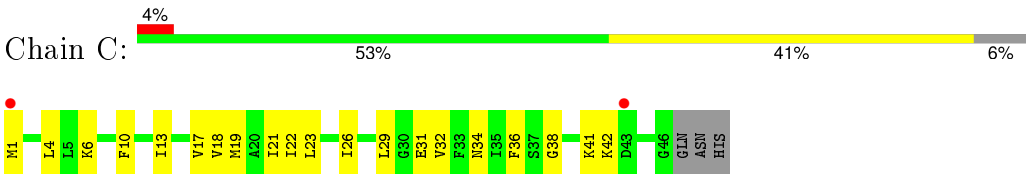
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1032	Total	C	N	O	S	0	0	0
			7788	5013	1285	1446	44			

- Molecule 2 is a protein called PREDICTED PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	46	Total	C	N	O	S	0	0	0
			326	219	48	56	3			



● Molecule 2: PREDICTED PROTEIN



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	146.21Å 146.21Å 543.21Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	24.91 – 3.70 24.89 – 3.70	Depositor EDS
% Data completeness (in resolution range)	94.0 (24.91-3.70) 94.5 (24.89-3.70)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.24 (at 3.74Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, $R_{free}$	0.339 , 0.361 0.335 , 0.353	Depositor DCC
$R_{free}$ test set	1158 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	142.3	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.26 , 53.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 23073 reflections	Xtriage
$F_o, F_c$ correlation	0.87	EDS
Total number of atoms	8114	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	119.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality [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	0/7936	0.69	1/10780 (0.0%)
2	C	0.49	0/330	0.64	0/448
All	All	0.54	0/8266	0.69	1/11228 (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	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	449	LEU	CA-CB-CG	5.09	127.00	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	222	THR	Mainchain,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	7788	0	7885	279	0
2	C	326	0	346	23	0
All	All	8114	0	8231	295	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 18.

All (295) 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)
2:C:6:LYS:HG3	2:C:10:PHE:CE2	1.53	1.43
2:C:6:LYS:CG	2:C:10:PHE:HE2	1.52	1.20
2:C:6:LYS:CG	2:C:10:PHE:CE2	2.35	0.96
1:A:575:MET:CE	1:A:662:MET:CE	2.49	0.90
1:A:70:ASN:O	1:A:110:LYS:NZ	2.08	0.86
1:A:555:LEU:HD22	1:A:913:LEU:HB3	1.59	0.85
1:A:960:LEU:HD21	1:A:1030:ARG:HB2	1.58	0.84
2:C:6:LYS:HG3	2:C:10:PHE:HE2	0.69	0.83
2:C:6:LYS:O	2:C:10:PHE:HD2	1.59	0.83
1:A:575:MET:HE2	1:A:662:MET:CE	2.11	0.80
1:A:540:ARG:HD2	2:C:38:GLY:HA3	1.64	0.79
1:A:291:ILE:HD13	1:A:306:ILE:HD13	1.66	0.78
2:C:6:LYS:O	2:C:10:PHE:CD2	2.37	0.78
1:A:575:MET:HE1	1:A:662:MET:CE	2.15	0.77
1:A:575:MET:CE	1:A:662:MET:HE2	2.14	0.77
1:A:575:MET:CE	1:A:662:MET:HE3	2.15	0.76
1:A:575:MET:HE2	1:A:662:MET:HE2	1.70	0.73
1:A:375:VAL:HG11	1:A:405:LEU:HD22	1.73	0.70
1:A:568:ASP:OD1	1:A:634:TRP:NE1	2.24	0.70
1:A:540:ARG:HD2	2:C:38:GLY:CA	2.21	0.70
1:A:684:LEU:HD22	1:A:695:LEU:HD11	1.72	0.70
1:A:901:VAL:HG21	1:A:943:ILE:HG13	1.72	0.70
1:A:938:SER:HB3	1:A:1014:ALA:HB1	1.73	0.70
1:A:360:GLN:O	1:A:361:ASN:ND2	2.25	0.69
1:A:632:LYS:O	1:A:637:ARG:NH1	2.25	0.69
1:A:422:GLU:N	1:A:422:GLU:OE1	2.25	0.69
1:A:57:VAL:HG21	1:A:86:GLY:HA2	1.75	0.69
1:A:446:ALA:HB2	1:A:482:VAL:HG21	1.75	0.68
1:A:454:VAL:HG12	1:A:475:VAL:HG21	1.75	0.68
1:A:137:LEU:HD21	1:A:303:ALA:HB2	1.75	0.67
1:A:633:ASP:OD1	1:A:634:TRP:N	2.27	0.67
1:A:721:LEU:HD13	1:A:815:ARG:HD3	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:598:TYR:HE2	1:A:629:VAL:HG21	1.61	0.66
1:A:248:LYS:HA	1:A:261:LEU:HD23	1.79	0.64
1:A:1030:ARG:O	1:A:1034:SER:OG	2.14	0.64
1:A:898:PRO:O	1:A:901:VAL:HG12	1.98	0.64
1:A:899:PHE:O	1:A:903:LEU:HD12	1.98	0.64
1:A:6:ILE:O	1:A:428:LYS:NZ	2.30	0.63
1:A:451:ALA:HB1	1:A:883:VAL:HG13	1.81	0.63
1:A:546:LEU:O	1:A:549:VAL:HG12	1.99	0.63
1:A:38:ILE:HG22	1:A:462:SER:HB2	1.80	0.63
1:A:424:GLY:HA3	1:A:502:LYS:HB3	1.80	0.62
1:A:712:MET:SD	1:A:835:LYS:HE2	2.38	0.62
1:A:402:ILE:O	1:A:406:VAL:HG13	1.99	0.62
1:A:884:VAL:HG12	1:A:902:MET:CE	2.30	0.62
1:A:984:LEU:HA	1:A:987:MET:HG2	1.82	0.62
1:A:453:PHE:HB3	1:A:932:LEU:HD22	1.80	0.62
1:A:453:PHE:CE1	1:A:475:VAL:HG22	2.35	0.61
1:A:485:ALA:O	1:A:490:PRO:HD3	2.00	0.61
1:A:110:LYS:HA	1:A:113:LEU:HG	1.83	0.60
1:A:926:TYR:HB3	1:A:1003:VAL:HG23	1.83	0.60
1:A:104:GLN:O	1:A:108:GLN:HG2	2.02	0.60
1:A:37:THR:HG21	1:A:296:GLY:HA2	1.82	0.60
1:A:903:LEU:O	1:A:906:PRO:HD2	2.02	0.60
1:A:20:MET:HA	1:A:377:LEU:HD11	1.82	0.60
1:A:404:LEU:HD11	1:A:449:LEU:HD21	1.83	0.59
1:A:153:ASP:OD1	1:A:182:TYR:OH	2.20	0.59
1:A:375:VAL:CG1	1:A:405:LEU:HD22	2.32	0.59
1:A:35:TYR:HE2	1:A:672:VAL:H	1.50	0.59
1:A:898:PRO:O	1:A:902:MET:HG2	2.03	0.59
1:A:885:PHE:HA	1:A:902:MET:HE3	1.84	0.59
1:A:679:GLY:HA2	1:A:830:GLN:HB3	1.85	0.58
1:A:196:PHE:O	1:A:252:LYS:NZ	2.27	0.58
1:A:448:VAL:CG2	1:A:884:VAL:HG13	2.34	0.58
1:A:637:ARG:HB2	1:A:642:ASN:HB3	1.85	0.58
1:A:1021:PHE:HB3	1:A:1025:PHE:CZ	2.39	0.57
1:A:36:PRO:O	1:A:38:ILE:HG13	2.04	0.57
1:A:901:VAL:HG23	1:A:942:ALA:HB3	1.86	0.57
1:A:94:PHE:HB3	1:A:98:THR:HG21	1.86	0.57
1:A:563:PHE:O	1:A:564:LEU:HD23	2.03	0.57
1:A:139:VAL:HG22	1:A:290:GLY:HA2	1.87	0.57
1:A:644:VAL:O	1:A:648:THR:HG23	2.04	0.57
1:A:454:VAL:N	1:A:455:PRO:HD2	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:544:LEU:HD12	1:A:547:ILE:HG12	1.87	0.56
1:A:110:LYS:HB2	1:A:113:LEU:HD12	1.87	0.56
1:A:712:MET:HG2	1:A:843:LEU:HD22	1.87	0.56
1:A:102:ILE:HA	1:A:105:VAL:HG22	1.87	0.56
2:C:13:ILE:O	2:C:17:VAL:HG23	2.06	0.56
1:A:425:LEU:HD12	1:A:429:GLU:HB3	1.87	0.56
1:A:169:THR:OG1	1:A:309:GLU:OE2	2.23	0.56
1:A:1032:ARG:NH2	2:C:42:LYS:O	2.39	0.56
1:A:653:ARG:O	1:A:656:SER:OG	2.25	0.55
1:A:575:MET:HE2	1:A:662:MET:HE3	1.79	0.55
1:A:396:PHE:O	1:A:399:VAL:HG12	2.05	0.55
1:A:568:ASP:OD2	1:A:644:VAL:HG23	2.07	0.55
1:A:888:LEU:HD12	1:A:902:MET:HE1	1.89	0.55
1:A:560:PRO:HG2	1:A:922:THR:HG22	1.89	0.55
1:A:83:ASP:OD1	1:A:87:THR:OG1	2.24	0.55
1:A:20:MET:HA	1:A:377:LEU:CD1	2.37	0.55
1:A:984:LEU:HD23	1:A:987:MET:HG3	1.90	0.54
1:A:330:THR:OG1	1:A:331:PRO:HD3	2.05	0.54
2:C:6:LYS:CD	2:C:10:PHE:HE2	2.20	0.54
1:A:339:GLU:O	1:A:343:THR:HG23	2.07	0.54
1:A:434:SER:HA	1:A:437:GLN:OE1	2.07	0.54
1:A:2:PRO:O	1:A:6:ILE:HD12	2.07	0.54
1:A:144:ASN:HA	1:A:320:GLY:O	2.08	0.54
1:A:719:ASN:HB2	1:A:828:LEU:HD13	1.89	0.54
1:A:476:SER:O	1:A:480:LEU:HG	2.08	0.53
1:A:449:LEU:HD13	1:A:936:GLY:HA3	1.89	0.53
1:A:137:LEU:HD13	1:A:293:LEU:HD21	1.90	0.53
1:A:412:VAL:O	1:A:416:VAL:HG23	2.08	0.53
1:A:973:ARG:NH1	1:A:977:MET:SD	2.81	0.53
1:A:721:LEU:HD13	1:A:815:ARG:CD	2.38	0.53
1:A:974:PRO:O	1:A:978:THR:HG23	2.08	0.53
1:A:588:GLN:O	1:A:592:ASN:ND2	2.41	0.53
1:A:415:ASN:ND2	1:A:437:GLN:HE21	2.05	0.53
1:A:358:PHE:CD2	1:A:977:MET:HG2	2.43	0.53
1:A:413:VAL:HA	1:A:493:CYS:SG	2.48	0.53
1:A:885:PHE:HA	1:A:902:MET:CE	2.39	0.53
1:A:1021:PHE:O	1:A:1024:VAL:HG22	2.09	0.53
1:A:464:GLY:O	1:A:468:ARG:HG2	2.09	0.53
1:A:477:ALA:O	1:A:481:SER:HB3	2.08	0.52
1:A:407:ASP:HA	1:A:410:ILE:HG12	1.89	0.52
1:A:971:ARG:NH1	1:A:974:PRO:HB2	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:932:LEU:O	1:A:935:ILE:HG13	2.09	0.52
1:A:398:MET:O	1:A:402:ILE:HG23	2.10	0.52
1:A:944:LEU:HD12	1:A:971:ARG:HH21	1.74	0.52
1:A:559:LEU:HD23	1:A:923:ASN:HB2	1.91	0.52
1:A:575:MET:HE1	1:A:662:MET:HE2	1.84	0.52
1:A:352:PHE:HD1	1:A:369:THR:HG21	1.73	0.52
1:A:449:LEU:HD23	1:A:478:MET:SD	2.50	0.52
1:A:905:VAL:HB	1:A:906:PRO:HD3	1.91	0.52
1:A:444:GLY:HA3	1:A:891:LEU:HD12	1.91	0.52
1:A:1032:ARG:NH1	2:C:41:LYS:CB	2.72	0.52
1:A:868:LEU:O	1:A:869:SER:OG	2.26	0.52
1:A:10:ILE:O	1:A:13:TRP:N	2.42	0.51
1:A:49:TYR:CD2	1:A:52:ALA:HB2	2.45	0.51
1:A:903:LEU:HD22	1:A:1025:PHE:CD1	2.46	0.51
1:A:417:GLU:O	1:A:420:MET:HG2	2.10	0.51
1:A:431:THR:HG21	1:A:494:ALA:HB2	1.93	0.51
1:A:904:VAL:HG11	1:A:1022:VAL:HG22	1.93	0.51
1:A:880:SER:O	1:A:884:VAL:HG23	2.11	0.51
1:A:26:ALA:O	1:A:30:LEU:HB2	2.10	0.51
1:A:997:SER:HA	1:A:1000:GLN:HE22	1.76	0.51
2:C:6:LYS:HG3	2:C:10:PHE:CD2	2.30	0.51
1:A:377:LEU:HD12	1:A:378:GLY:N	2.26	0.51
1:A:527:TYR:OH	1:A:1019:ILE:O	2.29	0.51
1:A:83:ASP:HA	1:A:815:ARG:HA	1.94	0.50
1:A:949:ALA:O	1:A:953:MET:HG2	2.11	0.50
1:A:1013:THR:O	1:A:1017:LEU:HB3	2.11	0.50
1:A:449:LEU:CD1	1:A:936:GLY:HA3	2.41	0.50
1:A:143:ILE:HD13	1:A:285:PRO:O	2.11	0.50
1:A:62:THR:O	1:A:66:GLU:HG3	2.11	0.50
1:A:35:TYR:OH	1:A:670:ALA:O	2.23	0.50
1:A:143:ILE:HD12	1:A:144:ASN:H	1.75	0.50
1:A:449:LEU:O	1:A:452:VAL:HG22	2.11	0.50
1:A:707:ALA:O	1:A:710:PRO:HD3	2.11	0.50
1:A:919:ARG:HH11	1:A:1005:THR:HG21	1.77	0.50
1:A:410:ILE:HD11	1:A:978:THR:HA	1.93	0.50
1:A:962:GLU:N	1:A:962:GLU:OE1	2.44	0.50
1:A:375:VAL:HG13	1:A:480:LEU:HD12	1.92	0.50
1:A:431:THR:O	1:A:435:MET:HG2	2.12	0.50
1:A:615:PHE:CD1	1:A:615:PHE:O	2.65	0.50
1:A:696:THR:HA	1:A:825:MET:HE1	1.95	0.49
1:A:453:PHE:HE1	1:A:475:VAL:HG22	1.75	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:989:LEU:HG	1:A:1000:GLN:HB2	1.93	0.49
1:A:135:SER:OG	1:A:672:VAL:O	2.17	0.49
1:A:198:LEU:HD13	1:A:251:LEU:HD13	1.93	0.49
1:A:102:ILE:O	1:A:106:GLN:HG3	2.13	0.49
1:A:448:VAL:O	1:A:452:VAL:HG13	2.13	0.49
1:A:369:THR:O	1:A:373:PRO:HD2	2.13	0.48
1:A:438:ILE:O	1:A:441:ALA:N	2.41	0.48
1:A:409:ALA:O	1:A:413:VAL:HG13	2.13	0.48
1:A:525:HIS:NE2	1:A:529:ASP:OD2	2.46	0.48
1:A:982:PHE:CD2	1:A:1011:MET:HG3	2.48	0.48
1:A:884:VAL:HG12	1:A:902:MET:HE1	1.95	0.48
1:A:649:MET:O	1:A:652:THR:OG1	2.26	0.48
1:A:913:LEU:O	1:A:917:THR:HG23	2.14	0.48
2:C:18:VAL:HA	2:C:21:ILE:HD12	1.94	0.48
1:A:686:ASP:HB2	1:A:695:LEU:HD22	1.96	0.48
2:C:6:LYS:CD	2:C:10:PHE:CE2	2.95	0.48
1:A:2:PRO:HB2	1:A:435:MET:HG3	1.96	0.48
1:A:719:ASN:ND2	1:A:826:GLU:OE2	2.47	0.48
1:A:440:GLY:O	1:A:891:LEU:HD13	2.14	0.48
1:A:437:GLN:HG2	1:A:438:ILE:HG23	1.96	0.48
1:A:940:LYS:HE3	1:A:941:ASN:OD1	2.14	0.48
1:A:595:THR:HG22	1:A:599:LEU:HD12	1.96	0.47
1:A:310:LEU:HD21	1:A:323:ILE:HD12	1.95	0.47
1:A:715:SER:O	1:A:829:GLY:HA2	2.14	0.47
1:A:445:ILE:HG22	1:A:943:ILE:HD13	1.97	0.47
1:A:584:GLN:HB2	1:A:622:GLN:HG2	1.96	0.47
1:A:415:ASN:O	1:A:419:VAL:HG23	2.15	0.47
1:A:583:THR:HG22	1:A:586:ARG:HD3	1.96	0.47
1:A:641:GLU:O	1:A:650:ARG:NH2	2.41	0.47
1:A:34:GLN:HB3	1:A:333:VAL:CG1	2.45	0.46
1:A:528:THR:HG21	1:A:969:ARG:HD3	1.96	0.46
1:A:901:VAL:HG23	1:A:942:ALA:CB	2.45	0.46
1:A:586:ARG:O	1:A:589:LYS:HB3	2.14	0.46
1:A:314:GLU:N	1:A:314:GLU:OE1	2.48	0.46
1:A:367:ILE:HB	1:A:368:PRO:HD3	1.98	0.46
1:A:151:GLN:HE22	1:A:278:ILE:HG23	1.80	0.46
1:A:980:LEU:HD22	2:C:19:MET:HE1	1.97	0.46
1:A:433:LYS:O	1:A:437:GLN:OE1	2.34	0.46
1:A:447:MET:O	1:A:450:SER:OG	2.27	0.46
1:A:471:SER:O	1:A:475:VAL:HG23	2.16	0.46
1:A:34:GLN:HB3	1:A:333:VAL:HG11	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:399:VAL:O	1:A:402:ILE:HG13	2.16	0.46
1:A:515:TRP:NE1	1:A:519:MET:SD	2.80	0.46
1:A:137:LEU:CD2	1:A:303:ALA:HB2	2.44	0.46
1:A:873:ALA:N	1:A:874:PRO:HD2	2.30	0.46
1:A:72:ILE:HG22	1:A:106:GLN:HB3	1.99	0.46
1:A:300:LEU:O	1:A:300:LEU:HD23	2.15	0.46
1:A:242:SER:OG	1:A:244:GLU:OE1	2.34	0.46
1:A:1:MET:HB3	1:A:2:PRO:HD3	1.97	0.45
1:A:643:LYS:O	1:A:647:ILE:HG13	2.15	0.45
1:A:211:ASN:HA	1:A:240:LEU:HD13	1.98	0.45
1:A:154:ILE:HG22	1:A:287:SER:HB3	1.99	0.45
1:A:782:LEU:O	1:A:785:ASP:HB2	2.17	0.45
1:A:960:LEU:HD21	1:A:1030:ARG:CB	2.37	0.45
1:A:987:MET:HB2	1:A:988:PRO:HD3	1.98	0.45
1:A:412:VAL:HB	1:A:438:ILE:HD11	1.97	0.45
1:A:722:GLU:O	1:A:724:THR:HG23	2.17	0.45
1:A:540:ARG:NH2	2:C:36:PHE:CB	2.80	0.45
1:A:1027:VAL:O	1:A:1031:ARG:N	2.45	0.45
1:A:559:LEU:HD12	1:A:560:PRO:HD2	2.00	0.45
1:A:110:LYS:HE2	1:A:110:LYS:HB3	1.55	0.44
1:A:888:LEU:HD13	1:A:901:VAL:HG11	2.00	0.44
1:A:176:GLN:N	1:A:290:GLY:O	2.50	0.44
1:A:754:TRP:CZ2	1:A:786:ILE:HD13	2.51	0.44
1:A:352:PHE:CD1	1:A:369:THR:HG21	2.52	0.44
1:A:184:MET:HB3	1:A:771:VAL:HG22	1.99	0.44
1:A:645:GLU:O	1:A:649:MET:HG2	2.18	0.44
1:A:711:ASP:N	1:A:711:ASP:OD1	2.51	0.44
1:A:358:PHE:CG	1:A:977:MET:HG2	2.53	0.44
2:C:23:LEU:HD23	2:C:26:ILE:HD11	1.99	0.44
1:A:402:ILE:HG13	1:A:403:GLY:N	2.33	0.44
1:A:16:ALA:O	1:A:20:MET:HG3	2.18	0.44
1:A:72:ILE:HG13	1:A:75:LEU:HD22	2.00	0.44
1:A:49:TYR:CE1	1:A:121:GLU:HB2	2.53	0.43
1:A:686:ASP:HB3	1:A:823:PRO:HB2	2.00	0.43
1:A:407:ASP:OD1	1:A:408:ASP:N	2.51	0.43
2:C:31:GLU:O	2:C:34:ASN:HB2	2.18	0.43
1:A:882:ILE:HD13	1:A:882:ILE:HA	1.94	0.43
1:A:832:ALA:HB1	1:A:833:PRO:HD2	2.00	0.43
1:A:102:ILE:O	1:A:105:VAL:HG22	2.19	0.43
1:A:594:VAL:HG13	1:A:655:PHE:CZ	2.52	0.43
2:C:29:LEU:HA	2:C:32:VAL:HG22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:LEU:HD11	1:A:376:LEU:HD22	2.00	0.43
1:A:480:LEU:O	1:A:484:VAL:HG13	2.18	0.43
1:A:453:PHE:CD1	1:A:475:VAL:HG22	2.53	0.43
1:A:988:PRO:O	1:A:991:ILE:HG13	2.19	0.43
1:A:958:LYS:HE3	1:A:962:GLU:HG2	2.00	0.43
1:A:428:LYS:HZ2	1:A:494:ALA:HB1	1.83	0.43
1:A:992:SER:CB	1:A:1000:GLN:HE21	2.32	0.43
1:A:111:LEU:HD12	1:A:129:VAL:HG23	2.00	0.42
1:A:535:LEU:HA	1:A:535:LEU:HD12	1.87	0.42
1:A:373:PRO:O	1:A:377:LEU:HG	2.18	0.42
1:A:45:ILE:HG12	1:A:129:VAL:HG22	2.01	0.42
1:A:552:MET:SD	1:A:909:VAL:HG21	2.60	0.42
1:A:403:GLY:O	1:A:406:VAL:HG22	2.19	0.42
1:A:699:ARG:HD3	1:A:825:MET:HE3	2.01	0.42
1:A:111:LEU:HD11	1:A:127:VAL:HG12	2.01	0.42
1:A:567:GLU:HG2	1:A:998:GLY:HA3	2.01	0.42
1:A:1008:MET:HB2	1:A:1008:MET:HE3	1.82	0.42
1:A:88:VAL:HG12	1:A:89:GLN:N	2.35	0.42
1:A:459:PHE:O	1:A:464:GLY:HA3	2.19	0.42
1:A:468:ARG:O	1:A:472:ILE:HG23	2.19	0.42
1:A:166:ILE:HG22	1:A:175:VAL:HG21	2.01	0.42
1:A:411:VAL:O	1:A:415:ASN:HB2	2.20	0.42
1:A:904:VAL:HA	1:A:907:LEU:HG	2.02	0.42
1:A:669:PRO:O	1:A:671:ILE:N	2.53	0.42
1:A:549:VAL:O	1:A:552:MET:HB3	2.19	0.41
2:C:19:MET:O	2:C:22:ILE:HG13	2.20	0.41
1:A:612:VAL:HG21	1:A:615:PHE:HB3	2.01	0.41
1:A:164:ASP:OD2	1:A:767:ARG:NH2	2.45	0.41
1:A:407:ASP:O	1:A:411:VAL:HG22	2.20	0.41
1:A:984:LEU:HA	1:A:987:MET:CG	2.49	0.41
1:A:944:LEU:HB3	1:A:971:ARG:HE	1.86	0.41
1:A:449:LEU:HG	1:A:453:PHE:CE2	2.55	0.41
1:A:928:GLN:O	1:A:932:LEU:HD12	2.21	0.41
1:A:895:TRP:O	1:A:896:SER:OG	2.28	0.41
1:A:644:VAL:HG11	1:A:667:ASN:ND2	2.36	0.41
1:A:448:VAL:HG11	1:A:888:LEU:HD21	2.02	0.41
1:A:1024:VAL:HA	1:A:1027:VAL:HG22	2.01	0.41
1:A:988:PRO:O	1:A:992:SER:HB2	2.21	0.41
1:A:166:ILE:HA	1:A:309:GLU:HG2	2.03	0.41
1:A:330:THR:O	1:A:333:VAL:N	2.53	0.41
1:A:186:ILE:HB	1:A:773:VAL:HG23	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:469:GLN:O	1:A:473:THR:OG1	2.28	0.41
1:A:980:LEU:HD21	2:C:19:MET:HB2	2.02	0.41
1:A:886:LEU:HD12	1:A:887:CYS:N	2.36	0.41
2:C:1:MET:O	2:C:4:LEU:HB2	2.21	0.40
1:A:399:VAL:HG21	1:A:989:LEU:HD11	2.03	0.40
1:A:49:TYR:HE1	1:A:121:GLU:HB2	1.86	0.40
1:A:342:LYS:O	1:A:346:GLU:HG2	2.20	0.40
1:A:891:LEU:HA	1:A:891:LEU:HD23	1.96	0.40
1:A:908:GLY:O	1:A:1010:GLY:HA2	2.20	0.40
1:A:454:VAL:N	1:A:455:PRO:CD	2.83	0.40
1:A:422:GLU:O	1:A:502:LYS:HB2	2.20	0.40
1:A:396:PHE:HE2	1:A:1003:VAL:HG21	1.87	0.40
1:A:1012:VAL:HG23	1:A:1013:THR:N	2.36	0.40
1:A:438:ILE:O	1:A:439:GLN:C	2.59	0.40
1:A:5:PHE:CD2	1:A:487:ILE:HG23	2.57	0.40
1:A:76:MET:SD	1:A:864:TYR:HE2	2.44	0.40
1:A:875:SER:O	1:A:878:ALA:HB3	2.21	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	1028/1049 (98%)	957 (93%)	68 (7%)	3 (0%)	46	83
2	C	44/49 (90%)	42 (96%)	2 (4%)	0	100	100
All	All	1072/1098 (98%)	999 (93%)	70 (6%)	3 (0%)	46	83

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	223	PRO

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Mol	Chain	Res	Type
1	A	224	PRO
1	A	579	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	823/855 (96%)	804 (98%)	19 (2%)	58	85
2	C	34/41 (83%)	34 (100%)	0	100	100
All	All	857/896 (96%)	838 (98%)	19 (2%)	60	86

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	SER
1	A	68	ASN
1	A	75	LEU
1	A	133	SER
1	A	412	VAL
1	A	425	LEU
1	A	472	ILE
1	A	481	SER
1	A	487	ILE
1	A	489	THR
1	A	493	CYS
1	A	547	ILE
1	A	672	VAL
1	A	801	PHE
1	A	904	VAL
1	A	932	LEU
1	A	945	ILE
1	A	993	THR
1	A	1034	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	360	GLN
1	A	361	ASN
1	A	415	ASN
1	A	517	ASN
1	A	592	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 carbohydrates in this entry.

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

There are no ligands in this entry.

### 5.7 Other polymers [i](#)

There are no such residues in this entry.

### 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	1032/1049 (98%)	-0.06	29 (2%) 56 41	104, 120, 120, 120	0
2	C	46/49 (93%)	-0.24	2 (4%) 39 26	120, 120, 120, 120	0
All	All	1078/1098 (98%)	-0.07	31 (2%) 55 39	104, 120, 120, 120	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1037	ASN	4.1
1	A	870	GLY	3.8
1	A	836	SER	3.6
1	A	1035	ARG	3.5
1	A	715	SER	3.2
1	A	1034	SER	3.2
1	A	829	GLY	2.9
1	A	679	GLY	2.8
1	A	109	ASN	2.8
1	A	833	PRO	2.8
1	A	717	ARG	2.7
1	A	173	GLY	2.7
1	A	869	SER	2.6
1	A	510	LYS	2.6
1	A	874	PRO	2.6
1	A	719	ASN	2.5
1	A	835	LYS	2.5
1	A	688	ALA	2.5
2	C	1	MET	2.4
1	A	840	ALA	2.3
1	A	720	GLY	2.3
2	C	43	ASP	2.3
1	A	690	LEU	2.2
1	A	461	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	638	PRO	2.2
1	A	834	GLY	2.2
1	A	98	THR	2.1
1	A	124	GLN	2.1
1	A	839	GLU	2.1
1	A	31	PRO	2.1
1	A	305	ALA	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 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.