



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:48 PM UTC

PDB ID : 5CSA / pdb_00005csa
Title : Crystal structure of domains BT-BCCP-AC1-AC5 of yeast acetyl-CoA carboxylase
Authors : Wei, J.; Tong, L.
Deposited on : 2015-07-23
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Xtrriage (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

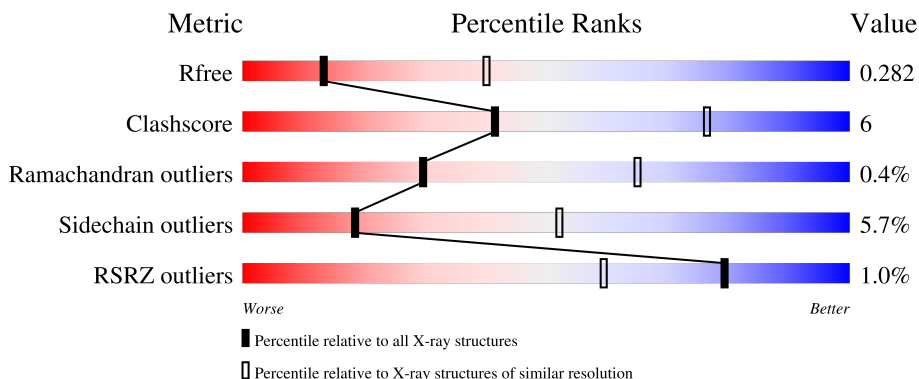
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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 ≥ 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	932	 % 75% 17% • 6%
1	B	932	 % 76% 14% • 9%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 13699 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 Acetyl-CoA carboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	872	6939	4440	1182	1296	21	0	4	0
1	B	852	6760	4329	1146	1265	20	0	0	0

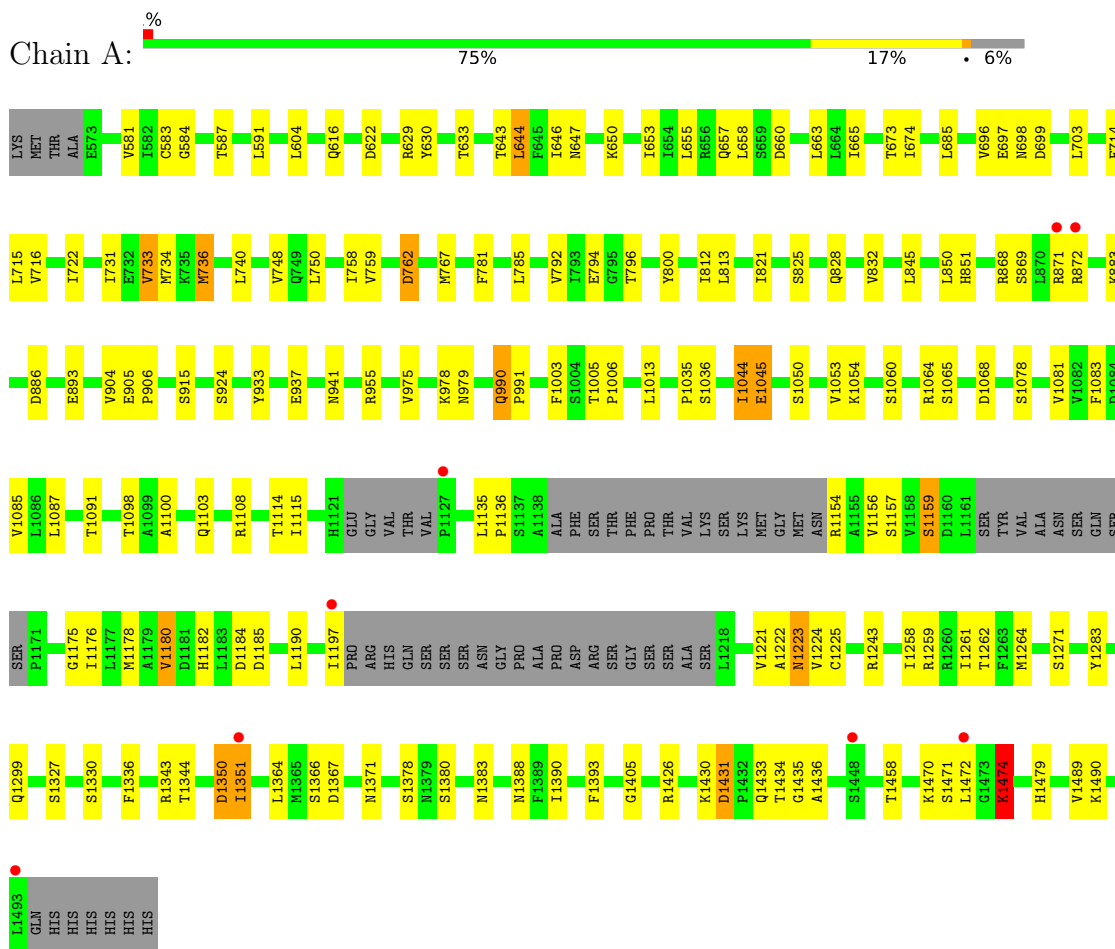
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1495	HIS	-	expression tag	UNP Q00955
A	1496	HIS	-	expression tag	UNP Q00955
A	1497	HIS	-	expression tag	UNP Q00955
A	1498	HIS	-	expression tag	UNP Q00955
A	1499	HIS	-	expression tag	UNP Q00955
A	1500	HIS	-	expression tag	UNP Q00955
B	1495	HIS	-	expression tag	UNP Q00955
B	1496	HIS	-	expression tag	UNP Q00955
B	1497	HIS	-	expression tag	UNP Q00955
B	1498	HIS	-	expression tag	UNP Q00955
B	1499	HIS	-	expression tag	UNP Q00955
B	1500	HIS	-	expression tag	UNP Q00955

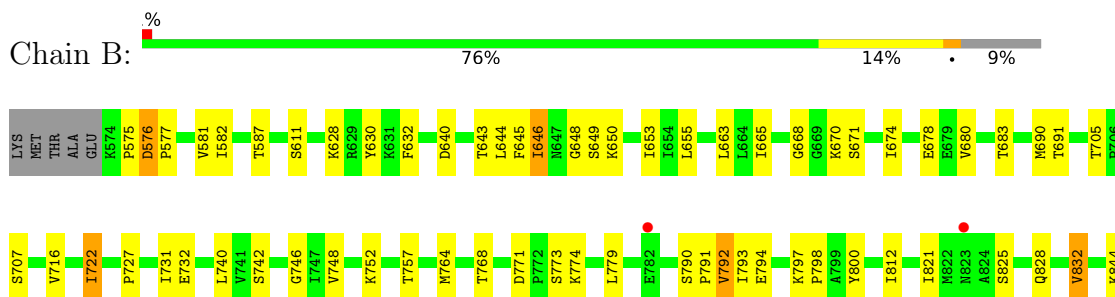
3 Residue-property plots i

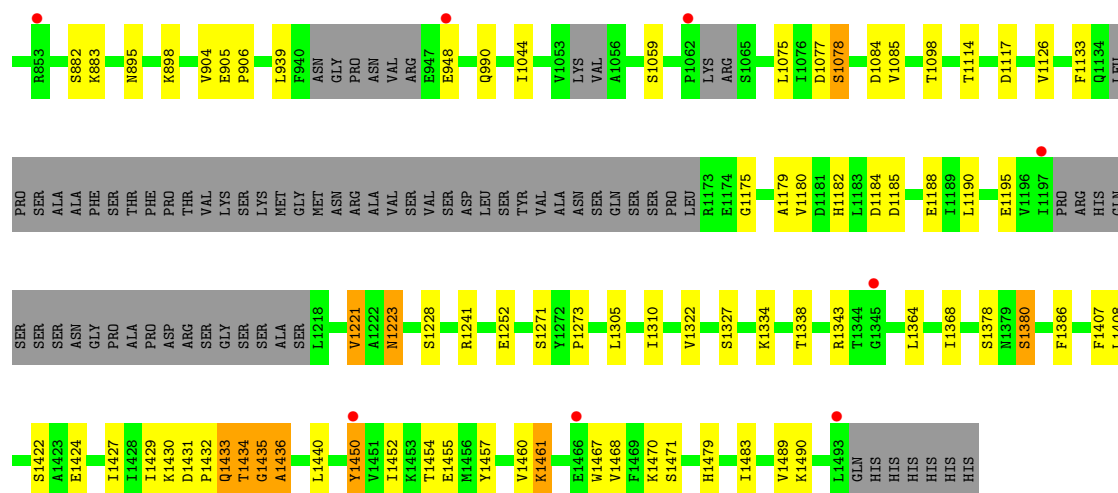
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Acetyl-CoA carboxylase



- Molecule 1: Acetyl-CoA carboxylase





4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	93.33Å 149.67Å 95.44Å 90.00° 118.39° 90.00°	Depositor
Resolution (Å)	42.63 – 3.00 42.63 – 3.00	Depositor EDS
% Data completeness (in resolution range)	97.8 (42.63-3.00) 97.8 (42.63-3.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.230 , 0.289 0.224 , 0.282	Depositor DCC
R_{free} test set	2290 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	60.3	Xtriage
Anisotropy	0.091	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 11.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for -h-l,k,h 0.000 for l,k,-h-l 0.020 for h,-k,-h-l 0.023 for -h-l,-k,l 0.018 for l,-k,h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	13699	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.73% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality i

5.1 Standard geometry i

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.72	0/7082	0.91	8/9586 (0.1%)
1	B	0.72	1/6890 (0.0%)	0.89	8/9327 (0.1%)
All	All	0.72	1/13972 (0.0%)	0.90	16/18913 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1450	TYR	CE1-CZ	7.00	1.55	1.38

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1133	PHE	N-CA-C	7.00	119.51	108.79
1	A	1433	GLN	N-CA-C	6.24	117.88	111.14
1	B	1435	GLY	N-CA-C	-6.02	107.35	115.36
1	B	1407	PHE	N-CA-C	-5.83	104.83	111.07
1	A	1405	GLY	N-CA-C	-5.60	105.37	112.54
1	A	1436	ALA	CA-C-N	-5.45	114.36	120.14
1	A	1436	ALA	C-N-CA	-5.45	114.36	120.14
1	B	1436	ALA	CA-C-N	-5.40	114.03	120.13
1	B	1436	ALA	C-N-CA	-5.40	114.03	120.13
1	A	1474	LYS	CD-CE-NZ	5.33	128.95	111.90
1	B	1378	SER	N-CA-C	5.24	116.69	110.97
1	B	576	ASP	CA-C-N	-5.08	113.49	119.84
1	B	576	ASP	C-N-CA	-5.08	113.49	119.84
1	A	1431	ASP	CA-C-N	-5.03	113.46	119.05
1	A	1431	ASP	C-N-CA	-5.03	113.46	119.05
1	A	1258	ILE	N-CA-C	5.01	114.82	107.15

There are no chirality outliers.

There are no planarity outliers.

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	6939	0	7077	94	6
1	B	6760	0	6877	89	6
All	All	13699	0	13954	178	6

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 6.

All (178) 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:1431:ASP:OD2	1:A:1434:THR:OG1	1.89	0.91
1:B:1461:LYS:HE2	1:B:1467:TRP:CD1	2.10	0.85
1:B:1432:PRO:HG2	1:B:1433:GLN:NE2	1.92	0.85
1:B:1432:PRO:HG2	1:B:1433:GLN:HE21	1.42	0.85
1:B:1431:ASP:OD2	1:B:1434:THR:OG1	2.00	0.79
1:A:872:ARG:NE	1:B:1188:GLU:HG2	2.00	0.77
1:B:1471:SER:HB2	1:B:1479:HIS:HD2	1.50	0.75
1:A:1154:ARG:NH2	1:A:1351:ILE:O	2.20	0.74
1:B:1327:SER:OG	1:B:1489:VAL:HG11	1.89	0.71
1:A:872:ARG:HE	1:B:1188:GLU:HG2	1.56	0.71
1:A:1431:ASP:CG	1:A:1434:THR:HG1	1.95	0.71
1:A:1371[B]:ASN:O	1:A:1371[B]:ASN:OD1	2.10	0.69
1:A:1431:ASP:OD2	1:A:1434:THR:N	2.21	0.69
1:A:722:ILE:HD12	1:A:748:VAL:HG21	1.75	0.68
1:A:1053:VAL:HG12	1:A:1054:LYS:O	1.94	0.67
1:B:1433:GLN:HE21	1:B:1433:GLN:N	1.93	0.67
1:A:1371[B]:ASN:OD1	1:A:1371[B]:ASN:C	2.37	0.66
1:A:1243:ARG:NH1	1:A:1283:TYR:O	2.26	0.66
1:A:1197:ILE:HG22	1:A:1197:ILE:O	1.96	0.64
1:A:716:VAL:HG21	1:A:722:ILE:HD11	1.79	0.64
1:B:731:ILE:HD11	1:B:740:LEU:HD11	1.80	0.64
1:B:792:VAL:HG11	1:B:800:TYR:CE2	2.34	0.63
1:A:872:ARG:CD	1:B:1188:GLU:HG2	2.29	0.63
1:B:1434:THR:HB	1:B:1436:ALA:H	1.64	0.61
1:A:868:ARG:HG2	1:A:871:ARG:NH1	2.16	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1305:LEU:HB3	1:B:1310:ILE:HD11	1.83	0.61
1:B:1252:GLU:N	1:B:1252:GLU:OE1	2.32	0.61
1:B:828:GLN:O	1:B:832:VAL:HG23	2.01	0.59
1:A:1078:SER:O	1:A:1108:ARG:NH2	2.35	0.58
1:B:1180:VAL:HG13	1:B:1185:ASP:HB2	1.86	0.58
1:A:937:GLU:OE1	1:A:975:VAL:HG11	2.03	0.58
1:B:628:LYS:HA	1:B:779:LEU:O	2.04	0.58
1:A:1458:THR:OG1	1:A:1472:LEU:HD21	2.04	0.58
1:B:1327:SER:OG	1:B:1489:VAL:CG1	2.51	0.58
1:A:1135:LEU:HB3	1:A:1136:PRO:CD	2.34	0.57
1:A:1197:ILE:O	1:A:1197:ILE:CG2	2.53	0.57
1:A:731:ILE:HD12	1:A:740:LEU:HD21	1.85	0.57
1:A:828:GLN:O	1:A:832:VAL:HG23	2.05	0.57
1:B:792:VAL:HG11	1:B:800:TYR:HE2	1.70	0.57
1:B:1455:GLU:OE1	1:B:1457:TYR:OH	2.19	0.57
1:A:1471:SER:HB2	1:A:1479:HIS:HD2	1.69	0.56
1:A:792:VAL:HG13	1:A:800:TYR:CE2	2.40	0.56
1:B:722:ILE:HD12	1:B:748:VAL:HG21	1.87	0.56
1:B:1182:HIS:CD2	1:B:1184:ASP:H	2.23	0.56
1:A:604:LEU:HG	1:A:955:ARG:HD3	1.88	0.56
1:A:587:THR:HG22	1:A:663:LEU:HD12	1.88	0.56
1:A:872:ARG:HD2	1:B:1188:GLU:HG2	1.88	0.56
1:A:698:ASN:OD1	1:A:699:ASP:N	2.39	0.55
1:A:1223:ASN:HB3	1:A:1262:THR:HB	1.89	0.55
1:A:655:LEU:HD22	1:A:663:LEU:HD22	1.89	0.55
1:A:583:CYS:SG	1:A:674:ILE:HD11	2.47	0.54
1:A:1225:CYS:HB2	1:A:1264:MET:HE3	1.88	0.54
1:A:658:LEU:O	1:A:660:ASP:O	2.25	0.54
1:B:794:GLU:O	1:B:844:LYS:CD	2.55	0.54
1:B:632:PHE:CZ	1:B:646:ILE:HD11	2.43	0.53
1:A:792:VAL:HG12	1:A:796:THR:OG1	2.10	0.52
1:B:1126:VAL:HG11	1:B:1179:ALA:O	2.08	0.52
1:B:771:ASP:OD1	1:B:773:SER:OG	2.28	0.52
1:A:1350:ASP:OD1	1:A:1350:ASP:N	2.43	0.52
1:B:731:ILE:HG22	1:B:732:GLU:N	2.25	0.52
1:B:1460:VAL:HG21	1:B:1470:LYS:HD3	1.91	0.51
1:A:587:THR:HG22	1:A:663:LEU:CD1	2.40	0.51
1:B:587:THR:HG22	1:B:663:LEU:HD12	1.93	0.51
1:A:715:LEU:O	1:A:716:VAL:HG23	2.11	0.50
1:A:714:PHE:CD1	1:A:750:LEU:HD22	2.46	0.50
1:B:752:LYS:HG2	1:B:764:MET:HA	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:941:ASN:HD21	1:A:1013:LEU:HA	1.75	0.50
1:B:1322:VAL:HG22	1:B:1338:THR:HG23	1.93	0.50
1:A:1064:ARG:HD2	1:A:1068:ASP:HB2	1.94	0.50
1:A:734:MET:C	1:A:736:MET:H	2.20	0.49
1:B:1431:ASP:O	1:B:1435:GLY:N	2.41	0.49
1:B:794:GLU:O	1:B:844:LYS:HD2	2.11	0.49
1:B:1432:PRO:CG	1:B:1433:GLN:HE21	2.19	0.49
1:B:1430:LYS:HE3	1:B:1435:GLY:HA2	1.94	0.49
1:B:665:ILE:HD12	1:B:674:ILE:CD1	2.43	0.49
1:B:1467:TRP:HB3	1:B:1483:ILE:HD12	1.94	0.49
1:B:1334:LYS:O	1:B:1380:SER:HA	2.12	0.49
1:A:584:GLY:HA2	1:A:685:LEU:HD11	1.95	0.49
1:B:895:ASN:OD1	1:B:898:LYS:HA	2.12	0.49
1:A:630:TYR:CE1	1:A:781:PHE:CD1	3.01	0.49
1:A:905:GLU:HB3	1:A:906:PRO:HD3	1.95	0.48
1:B:794:GLU:O	1:B:844:LYS:HD3	2.14	0.48
1:A:1115:ILE:HD12	1:A:1159:SER:HB3	1.94	0.47
1:A:644:LEU:O	1:A:650:LYS:HA	2.14	0.47
1:A:759:VAL:O	1:A:762:ASP:HB2	2.15	0.47
1:B:1195:GLU:O	1:B:1195:GLU:CD	2.57	0.47
1:B:1433:GLN:NE2	1:B:1433:GLN:N	2.60	0.47
1:B:1180:VAL:CG1	1:B:1185:ASP:HB2	2.44	0.47
1:B:1077:ASP:O	1:B:1078:SER:C	2.57	0.47
1:B:793:ILE:HG22	1:B:794:GLU:N	2.30	0.47
1:A:1156:VAL:HG12	1:A:1157:SER:O	2.15	0.47
1:B:792:VAL:HG21	1:B:800:TYR:CE2	2.50	0.47
1:B:1182:HIS:HD2	1:B:1184:ASP:H	1.62	0.47
1:A:990:GLN:HG3	1:A:991:PRO:N	2.30	0.47
1:B:1175:GLY:HA2	1:B:1221:VAL:O	2.15	0.47
1:B:1432:PRO:CG	1:B:1433:GLN:NE2	2.73	0.46
1:A:734:MET:O	1:A:736:MET:N	2.47	0.46
1:A:1044:ILE:HD12	1:A:1085:VAL:HG23	1.97	0.46
1:B:705:THR:HG22	1:B:740:LEU:HD11	1.97	0.46
1:A:1489:VAL:HG12	1:A:1490:LYS:N	2.31	0.46
1:B:792:VAL:C	1:B:793:ILE:HG13	2.41	0.46
1:A:1221:VAL:HG12	1:A:1222:ALA:N	2.31	0.46
1:A:1182:HIS:CD2	1:A:1184:ASP:H	2.34	0.46
1:A:933:TYR:OH	1:A:979:ASN:ND2	2.49	0.46
1:B:1433:GLN:HE21	1:B:1433:GLN:H	1.60	0.46
1:B:731:ILE:CD1	1:B:740:LEU:HD21	2.46	0.45
1:B:1429:ILE:HD12	1:B:1440:LEU:CD1	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1135:LEU:HB3	1:A:1136:PRO:HD3	1.98	0.45
1:A:665:ILE:HG13	1:A:674:ILE:HD12	1.98	0.45
1:A:587:THR:HG21	1:A:674:ILE:HG21	1.99	0.45
1:B:582:ILE:CG2	1:B:653:ILE:HD11	2.46	0.45
1:B:646:ILE:C	1:B:648:GLY:H	2.24	0.45
1:A:731:ILE:HD11	1:A:740:LEU:HD11	1.99	0.45
1:B:668:GLY:C	1:B:670:LYS:H	2.25	0.45
1:A:1336:PHE:N	1:A:1383:ASN:OD1	2.47	0.44
1:A:1431:ASP:O	1:A:1435:GLY:N	2.47	0.44
1:A:1035:PRO:HB3	1:B:1059:SER:O	2.17	0.44
1:B:821:ILE:CG2	1:B:825:SER:HB3	2.47	0.44
1:B:1489:VAL:HG12	1:B:1490:LYS:N	2.33	0.44
1:A:1003:PHE:C	1:A:1006:PRO:HD2	2.42	0.44
1:A:1178:MET:HE3	1:A:1224:VAL:HG22	2.00	0.44
1:A:1431:ASP:O	1:A:1435:GLY:HA2	2.16	0.44
1:B:581:VAL:HG21	1:B:630:TYR:CZ	2.53	0.44
1:B:746:GLY:HA3	1:B:768:THR:O	2.18	0.44
1:B:1044:ILE:HD12	1:B:1085:VAL:HG23	2.00	0.44
1:B:1431:ASP:OD1	1:B:1432:PRO:HD2	2.17	0.44
1:A:1100:ALA:O	1:A:1103:GLN:HB2	2.18	0.44
1:B:904:VAL:O	1:B:905:GLU:C	2.61	0.44
1:A:1175:GLY:HA2	1:A:1221:VAL:O	2.18	0.43
1:A:715:LEU:C	1:A:716:VAL:HG23	2.44	0.43
1:A:813:LEU:HA	1:A:978:LYS:HG2	2.00	0.43
1:A:733:VAL:N	1:A:736:MET:O	2.45	0.43
1:A:1366:SER:O	1:A:1367:ASP:C	2.61	0.43
1:A:904:VAL:O	1:A:905:GLU:C	2.61	0.43
1:A:646:ILE:HG23	1:A:647:ASN:N	2.33	0.43
1:A:1344:THR:HG21	1:A:1393:PHE:HZ	1.84	0.43
1:B:792:VAL:HG11	1:B:800:TYR:OH	2.18	0.43
1:B:1386:PHE:HA	1:B:1424:GLU:O	2.18	0.43
1:A:665:ILE:HD12	1:A:674:ILE:CD1	2.50	0.42
1:A:1327:SER:OG	1:A:1489:VAL:HG11	2.19	0.42
1:A:699:ASP:OD1	1:A:699:ASP:C	2.62	0.42
1:A:812:ILE:HD11	1:A:821:ILE:HD11	2.01	0.42
1:B:727:PRO:HA	1:B:740:LEU:O	2.20	0.42
1:A:1005:THR:HB	1:A:1006:PRO:HD3	2.02	0.42
1:B:628:LYS:HG3	1:B:630:TYR:CZ	2.53	0.42
1:B:678:GLU:HG2	1:B:683:THR:HG22	2.02	0.42
1:B:797:LYS:HB2	1:B:798:PRO:HD3	2.00	0.42
1:B:1408:LEU:HD22	1:B:1452:ILE:HD13	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1083:PHE:O	1:A:1087:LEU:HG	2.20	0.42
1:B:790:SER:HB2	1:B:791:PRO:CD	2.50	0.42
1:A:850:LEU:O	1:A:851:HIS:C	2.63	0.42
1:B:1338:THR:HG21	1:B:1368:ILE:HG23	2.02	0.42
1:A:587:THR:O	1:A:591:LEU:HG	2.20	0.42
1:A:1180:VAL:HG13	1:A:1185:ASP:HB2	2.02	0.42
1:B:645:PHE:HA	1:B:649:SER:O	2.20	0.42
1:A:714:PHE:CG	1:A:750:LEU:HD22	2.55	0.41
1:A:792:VAL:HG13	1:A:800:TYR:HE2	1.84	0.41
1:A:1299:GLN:O	1:A:1388:ASN:ND2	2.46	0.41
1:B:1223:ASN:OD1	1:B:1223:ASN:N	2.53	0.41
1:A:1050:SER:HA	1:A:1053:VAL:HG23	2.02	0.41
1:B:812:ILE:HD11	1:B:821:ILE:HD11	2.03	0.41
1:A:696:VAL:O	1:A:697:GLU:C	2.63	0.41
1:B:771:ASP:OD2	1:B:774:LYS:HE3	2.20	0.41
1:A:646:ILE:HG21	1:A:785:LEU:HG	2.03	0.41
1:A:1044:ILE:O	1:A:1045:GLU:C	2.62	0.41
1:B:716:VAL:HG21	1:B:722:ILE:HD11	2.01	0.41
1:B:905:GLU:HB3	1:B:906:PRO:HD3	2.01	0.41
1:A:622:ASP:OD1	1:A:622:ASP:C	2.64	0.41
1:A:703:LEU:HD11	1:A:767:MET:HE3	2.03	0.41
1:A:1470:LYS:O	1:A:1471:SER:C	2.63	0.41
1:B:576:ASP:O	1:B:577:PRO:C	2.61	0.41
1:B:1431:ASP:O	1:B:1435:GLY:HA2	2.20	0.41
1:B:575:PRO:HB3	1:B:690:MET:HE3	2.03	0.41
1:B:632:PHE:CE1	1:B:646:ILE:CG1	3.04	0.41
1:B:1461:LYS:HE2	1:B:1467:TRP:CG	2.53	0.40
1:A:657:GLN:HA	1:A:663:LEU:HD23	2.03	0.40
1:B:1489:VAL:CG1	1:B:1490:LYS:N	2.84	0.40
1:B:1084:ASP:HA	1:B:1273:PRO:HD3	2.04	0.40

All (6) 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:1474:LYS:CD	1:B:1450:TYR:CE1[2_546]	1.72	0.48
1:A:1474:LYS:CE	1:B:1450:TYR:CE1[2_546]	1.85	0.35
1:A:1474:LYS:CE	1:B:1450:TYR:OH[2_546]	2.05	0.15
1:A:1474:LYS:CE	1:B:1450:TYR:CZ[2_546]	2.09	0.11
1:A:1474:LYS:NZ	1:B:1450:TYR:CE1[2_546]	2.10	0.10
1:A:1474:LYS:CD	1:B:1450:TYR:CZ[2_546]	2.17	0.03

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	866/932 (93%)	792 (92%)	73 (8%)	1 (0%)	48	80
1	B	840/932 (90%)	775 (92%)	60 (7%)	5 (1%)	21	56
All	All	1706/1864 (92%)	1567 (92%)	133 (8%)	6 (0%)	30	65

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	680	VAL
1	A	1378	SER
1	B	939	LEU
1	B	948	GLU
1	B	646	ILE
1	B	792	VAL

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	774/821 (94%)	725 (94%)	49 (6%)	16	48
1	B	752/821 (92%)	715 (95%)	37 (5%)	22	56
All	All	1526/1642 (93%)	1440 (94%)	86 (6%)	18	52

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	581	VAL
1	A	616	GLN
1	A	629	ARG
1	A	633	THR
1	A	643	THR
1	A	644	LEU
1	A	653	ILE
1	A	673	THR
1	A	733	VAL
1	A	736	MET
1	A	758	ILE
1	A	762	ASP
1	A	794	GLU
1	A	825	SER
1	A	845	LEU
1	A	869	SER
1	A	883	LYS
1	A	886	ASP
1	A	893	GLU
1	A	915	SER
1	A	924	SER
1	A	990	GLN
1	A	1036	SER
1	A	1044	ILE
1	A	1045	GLU
1	A	1060	SER
1	A	1065	SER
1	A	1081	VAL
1	A	1091	THR
1	A	1098	THR
1	A	1114	THR
1	A	1159	SER
1	A	1176	ILE
1	A	1180	VAL
1	A	1190	LEU
1	A	1223	ASN
1	A	1259	ARG
1	A	1261	ILE
1	A	1271	SER
1	A	1330	SER
1	A	1343	ARG
1	A	1350	ASP
1	A	1351	ILE

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Mol	Chain	Res	Type
1	A	1364	LEU
1	A	1380	SER
1	A	1390	ILE
1	A	1426	ARG
1	A	1430	LYS
1	A	1474	LYS
1	B	611	SER
1	B	640	ASP
1	B	643	THR
1	B	644	LEU
1	B	650	LYS
1	B	655	LEU
1	B	671	SER
1	B	691	THR
1	B	707	SER
1	B	722	ILE
1	B	742	SER
1	B	757	THR
1	B	832	VAL
1	B	882	SER
1	B	883	LYS
1	B	990	GLN
1	B	1075	LEU
1	B	1078	SER
1	B	1098	THR
1	B	1114	THR
1	B	1117	ASP
1	B	1190	LEU
1	B	1221	VAL
1	B	1223	ASN
1	B	1228	SER
1	B	1241	ARG
1	B	1271	SER
1	B	1343	ARG
1	B	1364	LEU
1	B	1380	SER
1	B	1422	SER
1	B	1427	ILE
1	B	1433	GLN
1	B	1434	THR
1	B	1454	THR
1	B	1461	LYS

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Mol	Chain	Res	Type
1	B	1468	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	721	HIS
1	A	818	ASN
1	A	846	HIS
1	A	979	ASN
1	A	1043	GLN
1	A	1182	HIS
1	A	1249	ASN
1	A	1279	ASN
1	A	1319	ASN
1	A	1433	GLN
1	A	1479	HIS
1	B	880	GLN
1	B	921	HIS
1	B	1009	HIS
1	B	1031	GLN
1	B	1182	HIS
1	B	1319	ASN
1	B	1371	ASN
1	B	1433	GLN
1	B	1446	ASN
1	B	1479	HIS

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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	872/932 (93%)	-0.15	8 (0%) 81 61	30, 71, 105, 142	4 (0%)
1	B	852/932 (91%)	-0.03	10 (1%) 76 55	44, 75, 111, 151	0
All	All	1724/1864 (92%)	-0.09	18 (1%) 79 59	30, 72, 109, 151	4 (0%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1493	LEU	3.7
1	A	1448	SER	3.2
1	A	1127	PRO	3.0
1	B	1197	ILE	2.6
1	B	1062	PRO	2.6
1	B	1450	TYR	2.5
1	A	872	ARG	2.4
1	B	782	GLU	2.3
1	A	1197	ILE	2.3
1	A	1493	LEU	2.2
1	B	1466	GLU	2.2
1	B	823	ASN	2.1
1	A	1351	ILE	2.1
1	B	1345	GLY	2.1
1	B	948	GLU	2.1
1	A	871	ARG	2.0
1	A	1472	LEU	2.0
1	B	853	ARG	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 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.