



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 12:03 pm GMT

PDB ID : 1LBE
Title : APLYSIA ADP RIBOSYL CYCLASE
Authors : Prasad, G.S.; Mcree, D.E.; Stura, E.A.; Levitt, D.G.; Lee, H.C.; Stout, C.D.
Deposited on : 1996-09-18
Resolution : 2.40 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

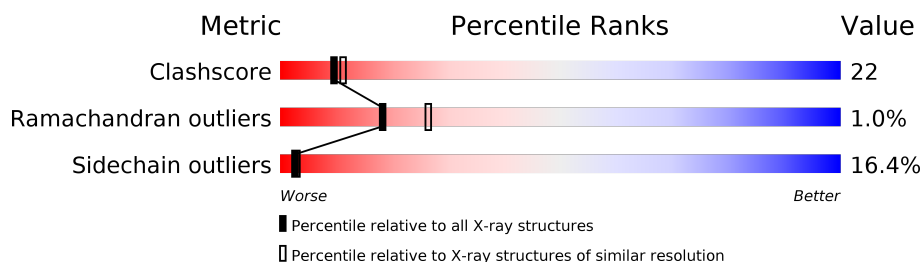
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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))
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)

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 ≥ 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	258	
1	B	258	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4064 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 ADP RIBOSYL CYCLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	250	Total	C	N	O	S	21	0	0
			2007	1285	341	367	14			
1	B	248	Total	C	N	O	S	13	0	0
			1988	1273	336	365	14			

- Molecule 2 is water.

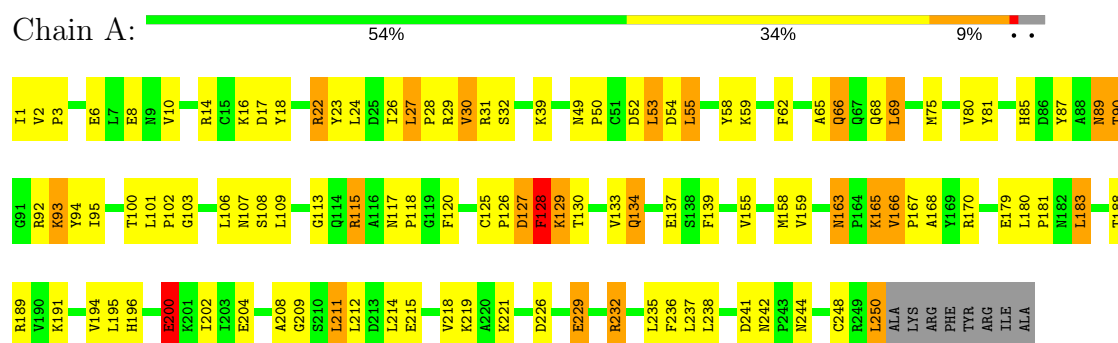
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	44	Total	O	0	0
			44	44		
2	B	25	Total	O	0	0
			25	25		

3 Residue-property plots [i](#)

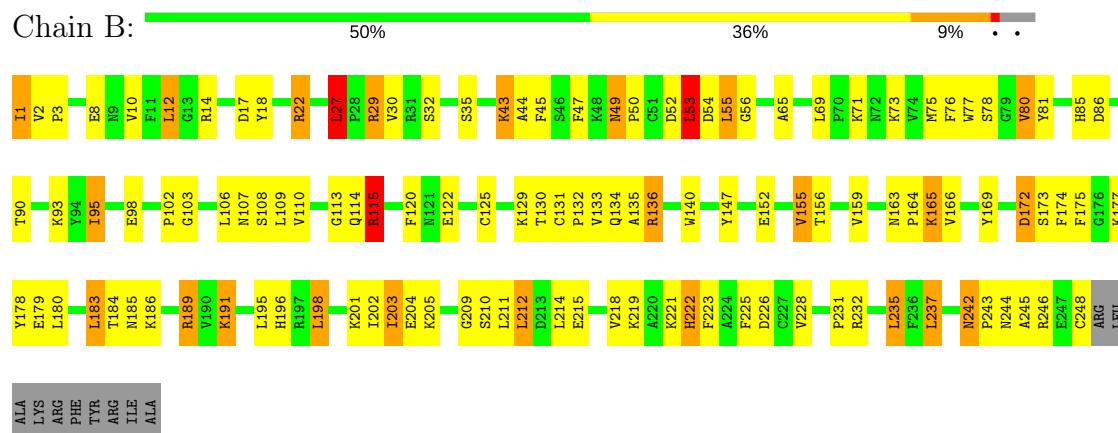
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

Note EDS was not executed.

• Molecule 1: ADP RIBOSYL CYCLASE



• Molecule 1: ADP RIBOSYL CYCLASE



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.20Å 72.30Å 125.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.40	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.40)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.214 , 0.299	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4064	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

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	0.88	0/2059	1.08	7/2786 (0.3%)
1	B	0.86	0/2040	1.04	7/2761 (0.3%)
All	All	0.87	0/4099	1.06	14/5547 (0.3%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	250	LEU	CA-CB-CG	9.90	138.06	115.30
1	B	106	LEU	CA-CB-CG	-7.75	97.48	115.30
1	B	29	ARG	NE-CZ-NH1	-6.89	116.86	120.30
1	A	29	ARG	NE-CZ-NH1	-6.81	116.89	120.30
1	B	53	LEU	CA-CB-CG	6.74	130.81	115.30
1	A	53	LEU	CA-CB-CG	6.47	130.18	115.30
1	A	22	ARG	N-CA-C	6.36	128.16	111.00
1	A	14	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	B	2	VAL	N-CA-C	-5.50	96.14	111.00
1	B	115	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	B	27	LEU	CA-CB-CG	5.39	127.70	115.30
1	A	200	GLU	N-CA-C	5.19	125.01	111.00
1	B	198	LEU	CA-CB-CG	5.07	126.97	115.30
1	A	208	ALA	C-N-CA	-5.06	111.68	122.30

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	2007	0	1963	78	0
1	B	1988	0	1939	97	0
2	A	44	0	0	6	0
2	B	25	0	0	3	0
All	All	4064	0	3902	169	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 22.

All (169) 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:175:PHE:HA	1:B:179:GLU:HB2	1.46	0.94
1:B:18:TYR:HA	1:B:22:ARG:HB2	1.60	0.84
1:B:184:THR:HG21	2:B:650:HOH:O	1.81	0.81
1:B:132:PRO:HB2	1:B:134:GLN:OE1	1.82	0.80
1:A:89:ASN:HB3	1:A:92:ARG:HB2	1.67	0.77
1:A:189:ARG:NH2	1:A:226:ASP:HB2	2.01	0.76
1:B:103:GLY:O	1:B:107:ASN:HB2	1.87	0.75
1:A:49:ASN:ND2	1:A:115:ARG:HH12	1.85	0.74
1:B:29:ARG:NH1	2:B:669:HOH:O	2.21	0.74
1:A:89:ASN:HB2	1:A:93:LYS:HE2	1.69	0.74
1:B:201:LYS:O	1:B:203:ILE:HD13	1.89	0.73
1:B:115:ARG:HG3	1:B:115:ARG:HH11	1.53	0.73
1:A:163:ASN:HD21	1:A:165:LYS:HG3	1.52	0.73
1:A:69:LEU:HG	1:A:95:ILE:HD11	1.70	0.72
1:B:173:SER:O	1:B:177:LYS:HG3	1.90	0.71
1:A:23:TYR:N	2:A:659:HOH:O	2.25	0.70
1:A:54:ASP:HA	1:A:134:GLN:HG2	1.74	0.69
1:B:80:VAL:HG11	1:B:195:LEU:HD22	1.75	0.68
1:A:26:ILE:H	1:A:26:ILE:HD12	1.57	0.68
1:B:202:ILE:HG21	1:B:205:LYS:HE2	1.75	0.68
1:B:133:VAL:HA	1:B:136:ARG:HB3	1.75	0.68
1:B:18:TYR:O	2:B:669:HOH:O	2.11	0.67
1:B:203:ILE:HG22	1:B:204:GLU:HG2	1.77	0.67
1:B:81:TYR:O	1:B:85:HIS:HD2	1.78	0.66
1:A:103:GLY:O	1:A:107:ASN:HB2	1.97	0.65
1:A:107:ASN:O	1:A:108:SER:HB2	1.96	0.65
1:B:69:LEU:HD21	1:B:147:TYR:CD1	2.32	0.65
1:B:1:ILE:HG13	1:B:125:CYS:HB2	1.78	0.64
1:B:221:LYS:O	1:B:222:HIS:HB2	1.97	0.64
1:B:215:GLU:HG2	1:B:219:LYS:HD3	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:PRO:HB3	1:B:125:CYS:SG	2.37	0.64
1:A:101:LEU:HB3	1:A:102:PRO:HD3	1.81	0.63
1:B:115:ARG:CG	1:B:115:ARG:HH11	2.12	0.62
1:B:1:ILE:HD12	1:B:110:VAL:CG1	2.30	0.61
1:B:201:LYS:HE3	1:B:203:ILE:HD12	1.82	0.61
1:B:65:ALA:HB1	1:B:102:PRO:HD3	1.83	0.60
1:B:198:LEU:HD11	1:B:248:CYS:SG	2.42	0.60
1:A:80:VAL:HG21	1:A:158:MET:HG2	1.83	0.59
1:A:236:PHE:HE2	1:B:243:PRO:HA	1.67	0.59
1:A:39:LYS:HD2	1:A:39:LYS:N	2.18	0.59
1:B:43:LYS:HG2	1:B:44:ALA:N	2.16	0.59
1:A:163:ASN:ND2	1:A:165:LYS:HG3	2.19	0.58
1:A:49:ASN:HB3	1:A:52:ASP:HB2	1.85	0.58
1:A:126:PRO:HB2	1:A:130:THR:HG21	1.85	0.58
1:A:89:ASN:O	1:A:92:ARG:HG3	2.02	0.58
1:A:90:THR:O	1:A:90:THR:CG2	2.52	0.58
1:B:242:ASN:ND2	1:B:244:ASN:O	2.36	0.57
1:A:163:ASN:C	1:A:163:ASN:HD22	2.08	0.57
1:A:49:ASN:OD1	1:A:50:PRO:HD2	2.05	0.57
1:B:77:TRP:CH2	1:B:81:TYR:HD2	2.22	0.57
1:B:1:ILE:HD12	1:B:110:VAL:HG11	1.86	0.56
1:B:78:SER:HB2	1:B:159:VAL:HG23	1.85	0.56
1:B:80:VAL:CG1	1:B:195:LEU:HD22	2.34	0.56
1:A:3:PRO:HB3	1:A:125:CYS:SG	2.45	0.56
1:A:165:LYS:HD2	1:A:166:VAL:HG12	1.88	0.56
1:A:106:LEU:HD13	1:A:139:PHE:CZ	2.41	0.55
1:B:164:PRO:HG2	1:B:165:LYS:H	1.71	0.55
1:A:179:GLU:O	1:A:183:LEU:HD13	2.05	0.54
1:A:18:TYR:HA	1:A:22:ARG:HG2	1.88	0.54
1:A:196:HIS:CE1	1:A:202:ILE:HG23	2.41	0.54
1:B:235:LEU:HD22	1:B:235:LEU:O	2.07	0.54
1:A:55:LEU:HD23	1:A:134:GLN:HB2	1.89	0.54
1:A:22:ARG:CA	2:A:659:HOH:O	2.56	0.54
1:A:133:VAL:O	1:A:137:GLU:N	2.42	0.53
1:B:214:LEU:O	1:B:218:VAL:HG23	2.08	0.53
1:A:17:ASP:OD1	1:B:14:ARG:HD2	2.08	0.53
1:B:8:GLU:O	1:B:12:LEU:HB2	2.08	0.53
1:B:76:PHE:CE2	1:B:155:VAL:HG13	2.44	0.53
1:A:163:ASN:ND2	1:A:165:LYS:H	2.06	0.52
1:B:53:LEU:O	1:B:135:ALA:HA	2.09	0.52
1:B:85:HIS:HE1	1:B:98:GLU:OE1	1.93	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:ASN:HD21	1:A:165:LYS:CG	2.21	0.52
1:B:169:TYR:CD1	1:B:214:LEU:HD13	2.45	0.52
1:A:248:CYS:O	1:B:232:ARG:NH2	2.42	0.52
1:B:172:ASP:O	1:B:177:LYS:HD2	2.10	0.52
1:A:163:ASN:OD1	1:A:166:VAL:HG13	2.10	0.51
1:A:90:THR:O	1:A:90:THR:HG23	2.10	0.51
1:A:163:ASN:HD22	1:A:165:LYS:H	1.57	0.51
1:A:232:ARG:NH1	1:B:243:PRO:O	2.41	0.51
1:B:218:VAL:CG1	1:B:223:PHE:HB2	2.41	0.51
1:A:39:LYS:H	1:A:39:LYS:HD2	1.76	0.51
1:B:228:VAL:CG1	1:B:231:PRO:HG3	2.41	0.51
1:A:89:ASN:O	1:A:90:THR:HB	2.11	0.49
1:B:174:PHE:CD2	1:B:178:TYR:HD2	2.30	0.49
1:B:189:ARG:HD2	1:B:226:ASP:HB2	1.94	0.49
1:B:196:HIS:CE1	1:B:204:GLU:O	2.65	0.49
1:A:18:TYR:HA	1:A:22:ARG:CG	2.42	0.49
1:A:236:PHE:CE2	1:B:243:PRO:HA	2.45	0.49
1:A:62:PHE:O	1:A:66:GLN:HB2	2.13	0.49
1:B:130:THR:HG22	1:B:131:CYS:O	2.12	0.49
1:A:209:GLY:N	2:A:668:HOH:O	2.44	0.49
1:B:133:VAL:HG12	1:B:136:ARG:HD2	1.94	0.49
1:B:140:TRP:C	1:B:174:PHE:HE2	2.16	0.49
1:A:65:ALA:HB1	1:A:102:PRO:HG3	1.94	0.49
1:B:169:TYR:CE1	1:B:214:LEU:HD13	2.48	0.48
1:B:49:ASN:HB3	1:B:115:ARG:NH2	2.28	0.48
1:B:76:PHE:HE2	1:B:155:VAL:CG1	2.26	0.48
1:A:159:VAL:HG21	1:A:168:ALA:O	2.13	0.48
1:B:196:HIS:HE1	1:B:204:GLU:O	1.95	0.48
1:B:76:PHE:HE2	1:B:155:VAL:HG13	1.79	0.48
1:A:167:PRO:HG2	1:A:170:ARG:HG3	1.96	0.47
1:B:218:VAL:HG21	1:B:225:PHE:CD1	2.49	0.47
1:A:30:VAL:HG13	1:A:32:SER:H	1.80	0.47
1:B:113:GLY:HA2	1:B:120:PHE:HA	1.97	0.47
1:B:10:VAL:CG1	1:B:109:LEU:HD21	2.45	0.46
1:A:221:LYS:NZ	2:A:631:HOH:O	2.46	0.46
1:A:128:PHE:HB2	1:A:129:LYS:H	1.63	0.46
1:B:245:ALA:HB3	1:B:248:CYS:SG	2.56	0.46
1:B:10:VAL:HG12	1:B:109:LEU:HD21	1.98	0.45
1:B:211:LEU:HD22	1:B:211:LEU:H	1.81	0.45
1:A:87:TYR:CD2	1:A:158:MET:HE1	2.50	0.45
1:B:81:TYR:O	1:B:85:HIS:CD2	2.64	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:ILE:HD12	1:B:110:VAL:HG13	1.96	0.45
1:A:26:ILE:N	1:A:26:ILE:HD12	2.29	0.45
1:B:177:LYS:HE2	1:B:177:LYS:HB3	1.74	0.45
1:B:47:PHE:HA	1:B:114:GLN:O	2.16	0.45
1:A:194:VAL:HB	1:A:229:GLU:HB3	1.98	0.45
1:A:214:LEU:O	1:A:218:VAL:HG23	2.17	0.45
1:B:75:MET:HA	1:B:156:THR:O	2.17	0.45
1:A:26:ILE:H	1:A:26:ILE:CD1	2.29	0.44
1:B:180:LEU:HA	1:B:183:LEU:HD22	1.98	0.44
1:B:242:ASN:N	1:B:243:PRO:HD3	2.33	0.44
1:A:75:MET:HE3	1:A:94:TYR:CD2	2.53	0.44
1:B:133:VAL:HA	1:B:136:ARG:HD2	2.00	0.44
1:B:228:VAL:HG11	1:B:231:PRO:HG3	2.00	0.44
1:B:49:ASN:O	1:B:52:ASP:HB2	2.17	0.44
1:A:118:PRO:HB2	1:A:120:PHE:CE1	2.53	0.44
1:B:49:ASN:HB2	1:B:115:ARG:NH1	2.33	0.44
1:A:196:HIS:HE1	1:A:204:GLU:O	2.01	0.43
1:A:211:LEU:HD13	1:A:211:LEU:HA	1.92	0.43
1:B:27:LEU:HD21	1:B:95:ILE:HD13	2.00	0.43
1:A:117:ASN:HD22	1:A:117:ASN:H	1.66	0.43
1:B:45:PHE:HA	1:B:53:LEU:HD11	2.01	0.42
1:B:169:TYR:HB3	1:B:210:SER:HB3	2.01	0.42
1:B:86:ASP:CB	1:B:237:LEU:HD11	2.50	0.42
1:A:58:TYR:HB2	1:A:62:PHE:HE2	1.85	0.42
1:B:54:ASP:O	1:B:56:GLY:N	2.53	0.42
1:A:1:ILE:HG23	1:A:125:CYS:HB2	2.01	0.42
1:A:24:LEU:HA	1:A:24:LEU:HD23	1.88	0.42
1:A:168:ALA:HB3	1:A:211:LEU:HD21	2.01	0.42
1:A:81:TYR:CE2	1:A:85:HIS:CE1	3.08	0.42
1:A:180:LEU:HB3	1:A:181:PRO:HD3	2.01	0.42
1:B:183:LEU:HA	1:B:183:LEU:HD12	1.90	0.42
1:A:27:LEU:HD12	1:A:28:PRO:HD3	2.01	0.42
1:B:50:PRO:HD3	1:B:113:GLY:O	2.19	0.42
1:B:131:CYS:HB3	1:B:132:PRO:HD2	2.01	0.42
1:A:244:ASN:HD22	1:B:232:ARG:NH1	2.18	0.42
1:B:73:LYS:HD3	1:B:93:LYS:O	2.20	0.42
1:A:22:ARG:N	2:A:659:HOH:O	2.52	0.41
1:B:1:ILE:N	1:B:125:CYS:O	2.47	0.41
1:A:50:PRO:HD3	1:A:113:GLY:O	2.20	0.41
1:A:120:PHE:N	1:A:120:PHE:CD1	2.87	0.41
1:A:215:GLU:O	1:A:219:LYS:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:THR:CB	2:A:617:HOH:O	2.68	0.41
1:A:6:GLU:O	1:A:10:VAL:HG23	2.20	0.41
1:B:107:ASN:O	1:B:108:SER:HB2	2.20	0.41
1:B:140:TRP:O	1:B:174:PHE:HE2	2.03	0.41
1:B:163:ASN:HA	1:B:164:PRO:HD2	1.91	0.41
1:B:55:LEU:HD23	1:B:134:GLN:HB2	2.02	0.41
1:B:12:LEU:HA	1:B:12:LEU:HD12	1.93	0.41
1:B:215:GLU:HG3	1:B:225:PHE:CD2	2.55	0.41
1:A:127:ASP:HB3	1:A:128:PHE:H	1.73	0.40
1:A:200:GLU:HG2	1:A:200:GLU:H	1.53	0.40
1:A:58:TYR:CB	1:A:62:PHE:CE2	3.04	0.40
1:B:212:LEU:O	1:B:212:LEU:HD22	2.21	0.40
1:B:221:LYS:HB3	1:B:223:PHE:CE2	2.56	0.40
1:B:69:LEU:HD21	1:B:147:TYR:CE1	2.55	0.40
1:B:71:LYS:HG3	1:B:152:GLU:HB2	2.03	0.40
1:B:191:LYS:HA	1:B:226:ASP:HB3	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	248/258 (96%)	233 (94%)	13 (5%)	2 (1%)	22	33
1	B	246/258 (95%)	223 (91%)	20 (8%)	3 (1%)	15	21
All	All	494/516 (96%)	456 (92%)	33 (7%)	5 (1%)	18	26

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	128	PHE
1	B	55	LEU

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Mol	Chain	Res	Type
1	B	222	HIS
1	B	209	GLY
1	A	127	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	220/226 (97%)	181 (82%)	39 (18%)	2	2
1	B	218/226 (96%)	185 (85%)	33 (15%)	3	4
All	All	438/452 (97%)	366 (84%)	72 (16%)	2	3

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	VAL
1	A	8	GLU
1	A	16	LYS
1	A	27	LEU
1	A	30	VAL
1	A	31	ARG
1	A	53	LEU
1	A	55	LEU
1	A	59	LYS
1	A	66	GLN
1	A	68	GLN
1	A	69	LEU
1	A	89	ASN
1	A	90	THR
1	A	93	LYS
1	A	109	LEU
1	A	115	ARG
1	A	128	PHE
1	A	129	LYS
1	A	134	GLN

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Mol	Chain	Res	Type
1	A	155	VAL
1	A	163	ASN
1	A	165	LYS
1	A	166	VAL
1	A	183	LEU
1	A	188	THR
1	A	191	LYS
1	A	195	LEU
1	A	200	GLU
1	A	211	LEU
1	A	212	LEU
1	A	229	GLU
1	A	232	ARG
1	A	235	LEU
1	A	237	LEU
1	A	238	LEU
1	A	241	ASP
1	A	242	ASN
1	A	250	LEU
1	B	1	ILE
1	B	12	LEU
1	B	17	ASP
1	B	22	ARG
1	B	27	LEU
1	B	30	VAL
1	B	32	SER
1	B	35	SER
1	B	43	LYS
1	B	49	ASN
1	B	53	LEU
1	B	80	VAL
1	B	90	THR
1	B	95	ILE
1	B	115	ARG
1	B	122	GLU
1	B	129	LYS
1	B	136	ARG
1	B	155	VAL
1	B	165	LYS
1	B	166	VAL
1	B	172	ASP
1	B	183	LEU

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Mol	Chain	Res	Type
1	B	185	ASN
1	B	186	LYS
1	B	189	ARG
1	B	191	LYS
1	B	203	ILE
1	B	212	LEU
1	B	235	LEU
1	B	237	LEU
1	B	242	ASN
1	B	246	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	GLN
1	A	117	ASN
1	A	163	ASN
1	A	196	HIS
1	A	242	ASN
1	A	244	ASN
1	B	85	HIS
1	B	196	HIS
1	B	242	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 ⓘ

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.