



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:04 PM GMT

PDB ID : 1NDB
Title : Crystal structure of Carnitine Acetyltransferase
Authors : Jogl, G.; Tong, L.
Deposited on : 2002-12-09
Resolution : 1.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
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
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

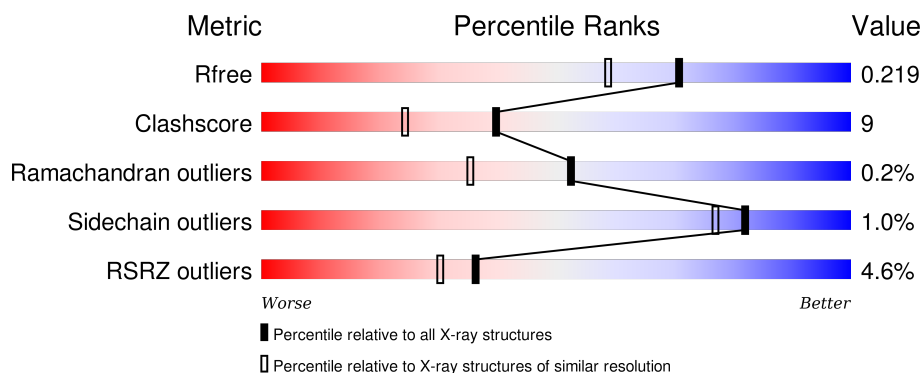
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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.

Mol	Chain	Length	Quality of chain
1	A	596	<div> <div>3%</div> <div>83%</div> <div>17%</div> </div>
1	B	596	<div> <div>6%</div> <div>80%</div> <div>19%</div> <div>.</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10560 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 Carnitine Acetyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	596	Total	C	N	O	S	Se	0	0	0
			4757	3034	820	875	8	20			
1	B	596	Total	C	N	O	S	Se	0	0	0
			4757	3034	820	875	8	20			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	94	MSE	MET	CLONING ARTIFACT	UNP P47934
A	152	MSE	MET	CLONING ARTIFACT	UNP P47934
A	170	MSE	MET	CLONING ARTIFACT	UNP P47934
A	308	MSE	MET	CLONING ARTIFACT	UNP P47934
A	339	MSE	MET	CLONING ARTIFACT	UNP P47934
A	359	MSE	MET	CLONING ARTIFACT	UNP P47934
A	372	MSE	MET	CLONING ARTIFACT	UNP P47934
A	377	MSE	MET	CLONING ARTIFACT	UNP P47934
A	403	MSE	MET	CLONING ARTIFACT	UNP P47934
A	410	MSE	MET	CLONING ARTIFACT	UNP P47934
A	411	MSE	MET	CLONING ARTIFACT	UNP P47934
A	459	MSE	MET	CLONING ARTIFACT	UNP P47934
A	482	MSE	MET	CLONING ARTIFACT	UNP P47934
A	534	MSE	MET	CLONING ARTIFACT	UNP P47934
A	539	MSE	MET	CLONING ARTIFACT	UNP P47934
A	547	MSE	MET	CLONING ARTIFACT	UNP P47934
A	564	MSE	MET	CLONING ARTIFACT	UNP P47934
A	581	MSE	MET	CLONING ARTIFACT	UNP P47934
A	603	MSE	MET	CLONING ARTIFACT	UNP P47934
A	614	MSE	MET	CLONING ARTIFACT	UNP P47934
B	94	MSE	MET	CLONING ARTIFACT	UNP P47934
B	152	MSE	MET	CLONING ARTIFACT	UNP P47934
B	170	MSE	MET	CLONING ARTIFACT	UNP P47934
B	308	MSE	MET	CLONING ARTIFACT	UNP P47934
B	339	MSE	MET	CLONING ARTIFACT	UNP P47934

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Chain	Residue	Modelled	Actual	Comment	Reference
B	359	MSE	MET	CLONING ARTIFACT	UNP P47934
B	372	MSE	MET	CLONING ARTIFACT	UNP P47934
B	377	MSE	MET	CLONING ARTIFACT	UNP P47934
B	403	MSE	MET	CLONING ARTIFACT	UNP P47934
B	410	MSE	MET	CLONING ARTIFACT	UNP P47934
B	411	MSE	MET	CLONING ARTIFACT	UNP P47934
B	459	MSE	MET	CLONING ARTIFACT	UNP P47934
B	482	MSE	MET	CLONING ARTIFACT	UNP P47934
B	534	MSE	MET	CLONING ARTIFACT	UNP P47934
B	539	MSE	MET	CLONING ARTIFACT	UNP P47934
B	547	MSE	MET	CLONING ARTIFACT	UNP P47934
B	564	MSE	MET	CLONING ARTIFACT	UNP P47934
B	581	MSE	MET	CLONING ARTIFACT	UNP P47934
B	603	MSE	MET	CLONING ARTIFACT	UNP P47934
B	614	MSE	MET	CLONING ARTIFACT	UNP P47934

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	584	Total O 584 584	0	0
2	B	462	Total O 462 462	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	158.88Å 89.65Å 119.44Å 90.00° 127.46° 90.00°	Depositor
Resolution (Å)	29.69 – 1.80 29.69 – 1.79	Depositor EDS
% Data completeness (in resolution range)	97.0 (29.69-1.80) 97.2 (29.69-1.79)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.50 (at 1.79Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.192 , 0.219 0.192 , 0.219	Depositor DCC
R_{free} test set	8795 reflections (7.29%)	DCC
Wilson B-factor (Å ²)	14.1	Xtriage
Anisotropy	0.224	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 57.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 240385 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10560	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.39% 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, 0.375 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	0.31	0/4852	0.57	0/6540
1	B	0.29	0/4852	0.56	0/6540
All	All	0.30	0/9704	0.56	0/13080

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4757	0	4733	86	0
1	B	4757	0	4733	80	0
2	A	584	0	0	12	0
2	B	462	0	0	6	0
All	All	10560	0	9466	166	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 9.

All (166) 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:372:MSE:HG2	1:A:373:VAL:H	1.31	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:LEU:HD22	1:A:132:LEU:H	1.39	0.87
1:A:624:ALA:O	1:A:625:LYS:HB2	1.76	0.86
1:B:623:ARG:HB3	1:B:623:ARG:NH1	1.90	0.85
1:A:32:GLN:HE22	1:A:170:MSE:H	1.28	0.81
1:B:265:ASN:HA	1:B:268:LYS:HE3	1.64	0.80
1:B:310:HIS:HD2	1:B:312:GLY:H	1.30	0.79
1:A:277:ASN:HD21	1:A:281:LYS:HE3	1.48	0.77
1:A:310:HIS:HD2	1:A:312:GLY:H	1.30	0.77
1:A:163:LEU:HB2	1:A:166:GLN:HG3	1.68	0.76
1:A:372:MSE:HG2	1:A:373:VAL:N	2.01	0.76
1:A:535:PRO:HB2	2:A:990:HOH:O	1.86	0.76
1:B:623:ARG:HB3	1:B:623:ARG:HH11	1.51	0.74
1:A:315:LYS:HG2	2:A:824:HOH:O	1.89	0.72
1:B:370:SER:HB2	1:B:371:PRO:HD2	1.73	0.69
1:A:310:HIS:CD2	1:A:312:GLY:H	2.12	0.68
1:B:349:PRO:HB2	1:B:350:PRO:HD3	1.76	0.67
1:A:459:MSE:HE3	1:A:511:ILE:HG22	1.75	0.67
1:A:255:HIS:HD2	1:A:257:ASN:H	1.41	0.67
1:A:537:ILE:HG22	2:A:990:HOH:O	1.94	0.66
1:B:117:TYR:CE1	1:B:403:MSE:HE3	2.30	0.66
1:B:132:LEU:HD12	1:B:133:GLN:H	1.60	0.66
1:B:310:HIS:CD2	1:B:312:GLY:H	2.12	0.66
1:A:349:PRO:HB2	1:A:350:PRO:HD3	1.78	0.66
1:A:123:ILE:HG13	1:A:565:PHE:CE2	2.32	0.65
1:A:322:TRP:H	1:A:328:GLN:NE2	1.96	0.64
1:A:568:PRO:HD2	1:A:592:TYR:CZ	2.33	0.64
1:A:277:ASN:ND2	1:A:281:LYS:HE3	2.12	0.64
1:A:255:HIS:CD2	1:A:257:ASN:H	2.15	0.63
1:A:512:ARG:HH11	1:A:512:ARG:HG3	1.64	0.63
1:A:132:LEU:CD2	1:A:132:LEU:H	2.12	0.61
1:A:358:VAL:HG12	1:A:359:MSE:HE2	1.82	0.61
1:B:421:PHE:HB3	1:B:422:PRO:HD3	1.82	0.61
1:A:132:LEU:N	1:A:132:LEU:HD22	2.16	0.59
1:B:497:ARG:O	1:B:501:GLN:HG2	2.03	0.59
1:B:95:GLU:OE1	1:B:461:HIS:HE1	1.86	0.59
1:A:322:TRP:H	1:A:328:GLN:HE22	1.51	0.58
1:B:104:LYS:HG3	1:B:105:THR:N	2.19	0.58
1:B:246:GLU:OE2	1:B:390:LYS:HE2	2.03	0.58
1:B:555:GLN:HE21	1:B:580:PRO:HG2	1.69	0.57
1:A:163:LEU:CB	1:A:166:GLN:HG3	2.34	0.57
1:A:555:GLN:HE21	1:A:580:PRO:HG2	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:276:VAL:O	1:B:280:GLN:HG3	2.04	0.56
1:A:493:VAL:HG22	1:A:617:LEU:HG	1.87	0.55
1:A:421:PHE:HB3	1:A:422:PRO:HD3	1.88	0.55
1:A:162:PHE:O	1:A:163:LEU:HD22	2.06	0.55
1:A:410:MSE:HG2	1:A:601:ALA:HA	1.88	0.55
1:B:457:LEU:HG	2:B:767:HOH:O	2.07	0.55
1:A:354:LEU:C	1:A:354:LEU:HD23	2.27	0.55
1:A:32:GLN:NE2	1:A:170:MSE:H	2.01	0.54
1:B:322:TRP:H	1:B:328:GLN:NE2	2.05	0.54
1:B:81:GLU:HG3	2:B:688:HOH:O	2.07	0.54
1:A:537:ILE:CG2	2:A:990:HOH:O	2.54	0.54
1:A:197:LYS:HG2	2:A:935:HOH:O	2.07	0.53
1:B:268:LYS:HG2	2:B:914:HOH:O	2.08	0.53
1:B:407:LEU:HD21	1:B:409:ILE:HD11	1.88	0.53
1:B:123:ILE:HG13	1:B:565:PHE:CE2	2.43	0.53
1:A:410:MSE:HE3	1:A:411:MSE:O	2.09	0.53
1:B:59:SER:OG	1:B:62:GLU:HG3	2.09	0.52
1:B:455:ALA:HB2	1:B:468:ILE:HG13	1.91	0.52
1:A:123:ILE:HD12	1:A:411:MSE:HE2	1.92	0.52
1:B:410:MSE:HE3	1:B:411:MSE:O	2.09	0.52
1:B:354:LEU:HD23	1:B:354:LEU:C	2.30	0.52
1:B:132:LEU:CD1	1:B:133:GLN:H	2.23	0.52
1:A:207:HIS:HD2	1:A:240:SER:OG	1.93	0.51
1:B:568:PRO:HD2	1:B:592:TYR:CZ	2.46	0.51
1:B:623:ARG:HH11	1:B:623:ARG:CB	2.22	0.51
1:A:163:LEU:CD2	1:A:168:LEU:HD11	2.40	0.51
1:B:129:PHE:CZ	1:B:335:GLY:HA2	2.45	0.51
1:A:129:PHE:CZ	1:A:335:GLY:HA2	2.45	0.51
1:B:555:GLN:HG2	1:B:557:PRO:HD3	1.92	0.51
1:B:416:HIS:CG	1:B:612:LEU:HD21	2.46	0.50
1:A:276:VAL:O	1:A:280:GLN:HG3	2.10	0.50
1:B:496:LEU:C	1:B:496:LEU:HD23	2.33	0.50
1:A:132:LEU:HD12	1:A:237:TRP:CZ2	2.47	0.49
1:B:116:ILE:HD12	1:B:403:MSE:HE2	1.93	0.49
1:B:81:GLU:O	1:B:85:LYS:HG2	2.12	0.49
1:A:231:VAL:HG21	1:A:372:MSE:HE1	1.93	0.49
1:A:163:LEU:HD23	1:A:168:LEU:HD11	1.95	0.49
1:B:202:HIS:HD2	2:B:666:HOH:O	1.95	0.49
1:A:421:PHE:CZ	1:A:615:ARG:HG3	2.48	0.48
1:A:123:ILE:HG13	1:A:565:PHE:CZ	2.48	0.48
1:A:457:LEU:HG	2:A:720:HOH:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:THR:HG21	1:A:279:ILE:HA	1.94	0.48
1:B:123:ILE:HD12	1:B:411:MSE:HE2	1.96	0.48
1:A:455:ALA:HB2	1:A:468:ILE:HG13	1.95	0.48
1:B:231:VAL:HG21	1:B:372:MSE:SE	2.62	0.48
1:B:255:HIS:HD2	1:B:257:ASN:HB2	1.78	0.48
1:B:322:TRP:H	1:B:328:GLN:HE22	1.60	0.48
1:A:294:ARG:HH11	1:A:294:ARG:HG3	1.78	0.48
1:B:349:PRO:HA	1:B:352:VAL:HG22	1.95	0.48
1:A:123:ILE:HG13	1:A:565:PHE:HE2	1.77	0.48
1:A:131:ASP:OD1	1:A:133:GLN:HG3	2.14	0.48
1:B:410:MSE:HG2	1:B:601:ALA:HA	1.96	0.47
1:B:204:THR:HG21	1:B:279:ILE:HA	1.96	0.47
1:A:265:ASN:HA	1:A:268:LYS:HE3	1.96	0.47
1:A:228:GLN:HG2	1:A:372:MSE:HG3	1.96	0.47
1:B:352:VAL:HG11	1:B:556:VAL:HG12	1.96	0.47
1:A:104:LYS:HG2	1:A:109:GLN:CD	2.35	0.47
1:B:115:VAL:HG12	1:B:116:ILE:HG13	1.96	0.47
1:A:198:ARG:HG3	1:A:198:ARG:O	2.15	0.47
1:A:401:SER:O	1:A:405:GLN:HG3	2.15	0.47
1:B:177:LEU:HD23	1:B:327:LEU:HD12	1.97	0.47
1:A:210:GLN:HE21	1:A:380:LYS:NZ	2.13	0.47
1:A:411:MSE:CE	1:A:563:VAL:HG11	2.45	0.46
1:A:115:VAL:HG12	1:A:116:ILE:HG13	1.98	0.46
1:A:141:LYS:HZ1	1:A:369:ARG:NH2	2.13	0.46
1:A:390:LYS:HB3	1:A:390:LYS:NZ	2.31	0.46
1:B:31:HIS:CE1	1:B:167:PRO:HB3	2.51	0.46
1:B:255:HIS:HE1	2:B:842:HOH:O	1.99	0.46
1:B:609:LYS:HE2	1:B:613:ASP:OD1	2.15	0.46
1:A:68:GLN:NE2	2:A:1063:HOH:O	2.49	0.46
1:A:548:HIS:HE1	1:A:572:ASP:OD1	1.99	0.45
1:A:405:GLN:HG3	2:A:953:HOH:O	2.15	0.45
1:B:202:HIS:HE1	1:B:278:SER:O	1.98	0.45
1:B:207:HIS:HD2	1:B:240:SER:OG	1.99	0.45
1:A:142:LEU:HD23	1:A:142:LEU:C	2.37	0.45
1:A:512:ARG:HG3	1:A:512:ARG:NH1	2.28	0.45
1:A:177:LEU:HD23	1:A:327:LEU:HD12	1.98	0.45
1:B:310:HIS:HE1	2:B:948:HOH:O	2.00	0.45
1:B:564:MSE:HB3	1:B:577:CYS:SG	2.57	0.45
1:A:255:HIS:CD2	1:A:257:ASN:HB2	2.51	0.45
1:A:310:HIS:HE1	2:A:1026:HOH:O	2.00	0.45
1:B:320:ASN:C	1:B:321:ARG:HG2	2.36	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:228:GLN:CG	1:B:372:MSE:HE2	2.47	0.44
1:A:496:LEU:C	1:A:496:LEU:HD23	2.37	0.44
1:B:548:HIS:HE1	1:B:572:ASP:OD1	1.99	0.44
1:B:228:GLN:HG3	1:B:372:MSE:HE2	1.99	0.44
1:B:68:GLN:O	1:B:72:GLU:HG3	2.18	0.44
1:B:124:LEU:HG	1:B:338:GLY:HA2	2.00	0.44
1:B:142:LEU:HD23	1:B:142:LEU:C	2.38	0.44
1:A:237:TRP:HH2	2:A:835:HOH:O	2.00	0.44
1:B:372:MSE:CG	1:B:373:VAL:N	2.79	0.44
1:B:123:ILE:HG13	1:B:565:PHE:CZ	2.53	0.44
1:B:204:THR:HG23	1:B:282:SER:HB3	1.99	0.44
1:A:349:PRO:HA	1:A:352:VAL:HG22	2.00	0.43
1:A:192:ASN:C	1:A:192:ASN:HD22	2.21	0.43
1:A:109:GLN:NE2	2:A:992:HOH:O	2.51	0.43
1:A:228:GLN:CG	1:A:372:MSE:HG3	2.48	0.43
1:B:50:TYR:CZ	1:B:54:LEU:HD11	2.54	0.43
1:A:202:HIS:HE1	1:A:278:SER:O	2.02	0.42
1:B:31:HIS:CE1	1:B:160:VAL:HG11	2.54	0.42
1:B:199:PRO:HA	1:B:200:PRO:HD3	1.89	0.42
1:A:95:GLU:OE1	1:A:461:HIS:HE1	2.02	0.42
1:B:192:ASN:C	1:B:192:ASN:HD22	2.22	0.42
1:B:157:THR:HG22	1:B:157:THR:O	2.19	0.42
1:B:176:ILE:HA	1:B:326:THR:HG21	2.01	0.42
1:A:294:ARG:HG3	1:A:294:ARG:NH1	2.35	0.42
1:B:65:HIS:CE1	1:B:69:LEU:HD11	2.54	0.42
1:A:564:MSE:HB3	1:A:577:CYS:SG	2.60	0.42
1:B:90:ARG:O	1:B:94:MSE:HG2	2.19	0.42
1:B:434:GLN:HG3	1:B:503:HIS:CE1	2.54	0.42
1:B:172:GLN:OE1	1:B:350:PRO:HG2	2.20	0.42
1:B:206:VAL:HG22	1:B:320:ASN:OD1	2.20	0.42
1:A:320:ASN:C	1:A:321:ARG:HG2	2.40	0.41
1:B:41:PRO:HA	1:B:42:PRO:HD3	1.95	0.41
1:A:199:PRO:HA	1:A:200:PRO:HD3	1.92	0.41
1:A:206:VAL:HG22	1:A:211:PHE:CD2	2.55	0.41
1:A:535:PRO:C	2:A:990:HOH:O	2.59	0.41
1:A:416:HIS:CG	1:A:612:LEU:HD21	2.55	0.41
1:B:487:VAL:HA	1:B:488:PRO:HD3	1.88	0.41
1:B:147:LEU:HD11	1:B:216:VAL:HB	2.02	0.41
1:B:349:PRO:O	1:B:352:VAL:HG22	2.21	0.41
1:B:192:ASN:ND2	1:B:194:LEU:H	2.19	0.41
1:B:35:LEU:HA	1:B:36:PRO:HD3	1.95	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:PHE:CE1	1:A:153:ILE:HD11	2.56	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	594/596 (100%)	578 (97%)	16 (3%)	0	100	100
1	B	594/596 (100%)	578 (97%)	14 (2%)	2 (0%)	46	29
All	All	1188/1192 (100%)	1156 (97%)	30 (2%)	2 (0%)	52	35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	624	ALA
1	B	623	ARG

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	524/504 (104%)	520 (99%)	4 (1%)	86	83
1	B	524/504 (104%)	517 (99%)	7 (1%)	76	68
All	All	1048/1008 (104%)	1037 (99%)	11 (1%)	82	77

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	166	GLN
1	A	192	ASN
1	A	343	HIS
1	A	506	TYR
1	B	132	LEU
1	B	156	GLU
1	B	166	GLN
1	B	192	ASN
1	B	405	GLN
1	B	506	TYR
1	B	615	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (53) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	65	HIS
1	A	68	GLN
1	A	84	GLN
1	A	109	GLN
1	A	112	GLN
1	A	135	GLN
1	A	166	GLN
1	A	187	GLN
1	A	192	ASN
1	A	202	HIS
1	A	207	HIS
1	A	210	GLN
1	A	228	GLN
1	A	255	HIS
1	A	277	ASN
1	A	307	GLN
1	A	310	HIS
1	A	328	GLN
1	A	357	HIS
1	A	461	HIS
1	A	503	HIS
1	A	548	HIS
1	A	550	ASN
1	A	555	GLN
1	A	586	ASN

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Mol	Chain	Res	Type
1	A	619	GLN
1	A	620	ASN
1	B	32	GLN
1	B	65	HIS
1	B	68	GLN
1	B	84	GLN
1	B	109	GLN
1	B	112	GLN
1	B	135	GLN
1	B	187	GLN
1	B	192	ASN
1	B	202	HIS
1	B	207	HIS
1	B	228	GLN
1	B	255	HIS
1	B	307	GLN
1	B	310	HIS
1	B	328	GLN
1	B	357	HIS
1	B	461	HIS
1	B	503	HIS
1	B	548	HIS
1	B	550	ASN
1	B	555	GLN
1	B	586	ASN
1	B	605	HIS
1	B	619	GLN

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 [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, 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	576/596 (96%)	-0.11	20 (3%)	48 42	5, 12, 29, 71	0
1	B	576/596 (96%)	0.12	33 (5%)	27 22	8, 17, 41, 70	0
All	All	1152/1192 (96%)	0.01	53 (4%)	36 30	5, 14, 33, 71	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	164	GLY	13.6
1	B	624	ALA	13.1
1	A	368	VAL	12.5
1	A	165	GLY	10.7
1	B	368	VAL	9.8
1	B	625	LYS	9.1
1	A	371	PRO	8.9
1	A	367	LEU	8.5
1	A	369	ARG	8.5
1	B	623	ARG	7.1
1	B	367	LEU	5.9
1	B	369	ARG	5.0
1	B	371	PRO	4.7
1	B	33	ASP	4.5
1	B	163	LEU	4.4
1	A	132	LEU	4.3
1	B	30	ALA	4.2
1	A	370	SER	4.2
1	B	165	GLY	4.0
1	B	32	GLN	3.9
1	A	166	GLN	3.8
1	A	163	LEU	3.7
1	B	370	SER	3.5
1	B	198	ARG	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	365	PRO	3.5
1	A	198	ARG	3.4
1	A	366	GLU	3.3
1	B	622	PRO	3.2
1	B	219	SER	3.2
1	A	33	ASP	3.0
1	B	373	VAL	3.0
1	B	132	LEU	3.0
1	B	164	GLY	3.0
1	B	34	ALA	2.9
1	B	31	HIS	2.7
1	B	157	THR	2.7
1	A	625	LYS	2.5
1	B	76	SER	2.4
1	B	374	PRO	2.4
1	B	364	LYS	2.4
1	A	365	PRO	2.4
1	A	162	PHE	2.4
1	A	30	ALA	2.3
1	A	237	TRP	2.3
1	B	156	GLU	2.3
1	A	294	ARG	2.3
1	A	243	SER	2.3
1	B	162	PHE	2.2
1	B	197	LYS	2.2
1	B	366	GLU	2.2
1	B	237	TRP	2.1
1	B	346	ALA	2.0
1	B	128	ASP	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.