



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:50 PM UTC

PDB ID : 1Q6X / pdb_00001q6x
Title : Crystal structure of rat choline acetyltransferase
Authors : Cai, Y.; Rodgers, D.W.
Deposited on : 2003-08-14
Resolution : 2.50 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
Xtriage (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

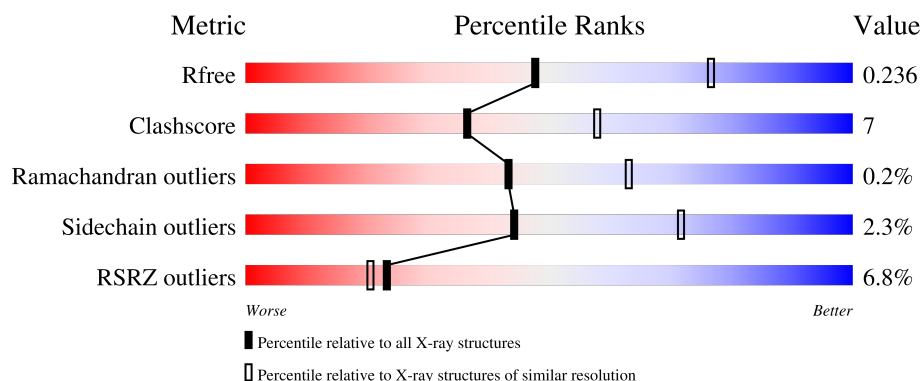
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 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	644	7% 77% 15% • 7%
1	B	644	6% 77% 14% • 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9681 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 choline O-acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	600	Total	C	N	O	S	0	0	12
			4667	2955	806	867	39			
1	B	600	Total	C	N	O	S	0	0	12
			4667	2955	806	867	39			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	234	ASN	LYS	SEE REMARK 999	UNP P32738
B	234	ASN	LYS	SEE REMARK 999	UNP P32738

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		
2	B	1	Total	Na	0	0
			1	1		

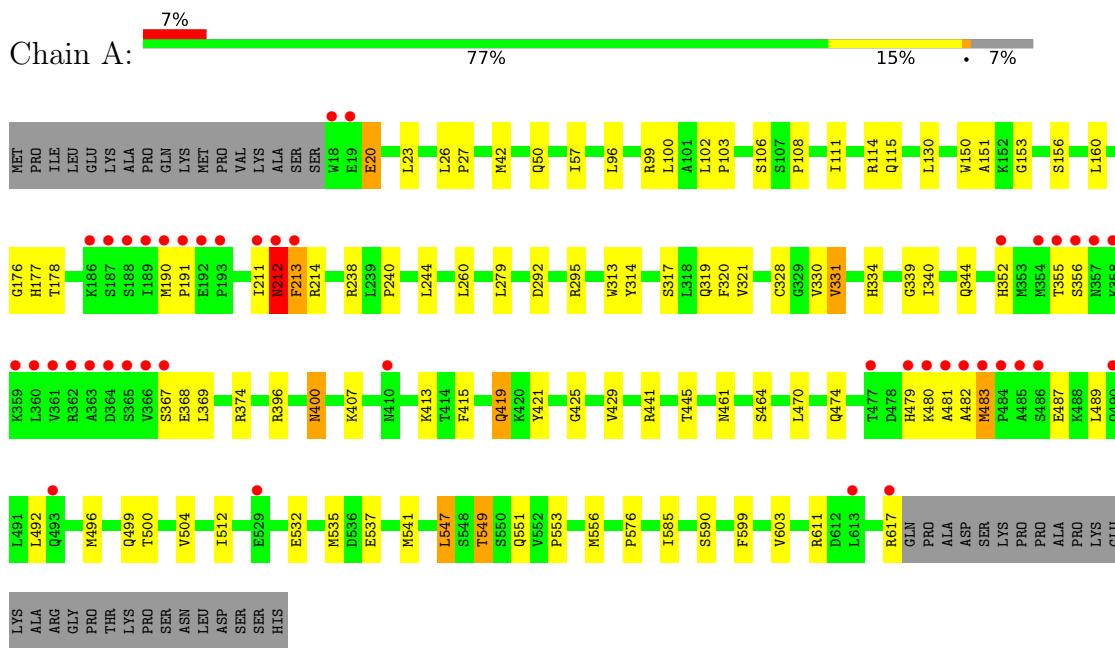
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	149	Total	O	0	0
			149	149		
3	B	196	Total	O	0	0
			196	196		

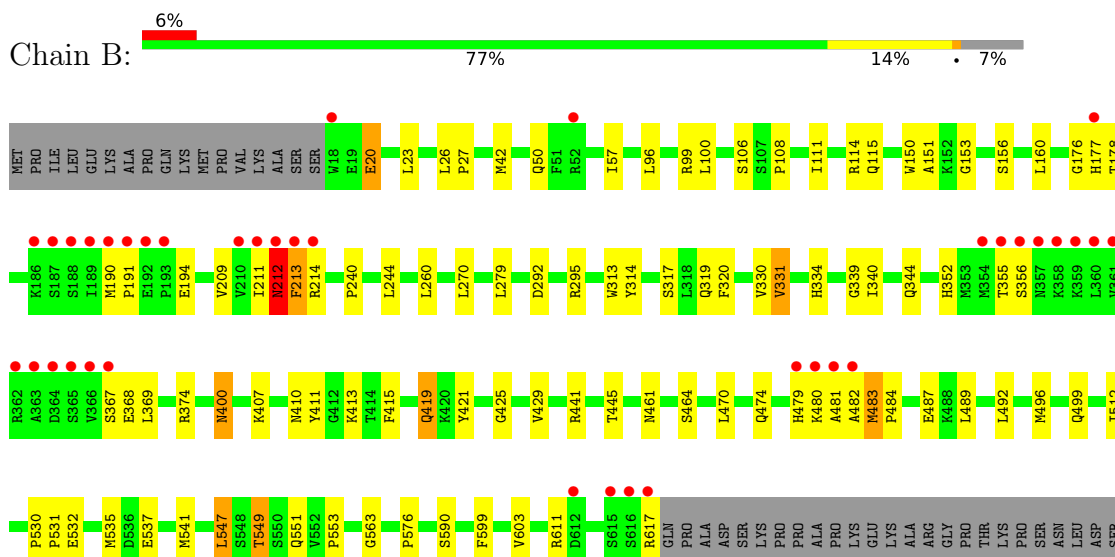
3 Residue-property plots

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: choline O-acetyltransferase



• Molecule 1: choline O-acetyltransferase



SER
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	84.70Å 78.88Å 139.36Å 90.00° 98.38° 90.00°	Depositor
Resolution (Å)	28.94 – 2.50 28.94 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.1 (28.94-2.50) 98.1 (28.94-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.93 (at 2.51Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.223 , 0.251 0.209 , 0.236	Depositor DCC
R_{free} test set	6279 reflections (10.12%)	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 31.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	9681	wwPDB-VP
Average B, all atoms (Å ²)	30.0	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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA

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	0.50	1/4756 (0.0%)	0.90	15/6442 (0.2%)
1	B	0.53	1/4756 (0.0%)	0.90	16/6442 (0.2%)
All	All	0.51	2/9512 (0.0%)	0.90	31/12884 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	240	PRO	CA-C	5.50	1.54	1.51
1	A	240	PRO	CA-C	5.17	1.54	1.51

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	314	TYR	N-CA-C	6.60	120.43	112.38
1	A	314	TYR	N-CA-C	6.56	120.38	112.38
1	A	213	PHE	N-CA-C	-6.48	105.34	112.72
1	B	213	PHE	N-CA-C	-6.34	105.49	112.72
1	A	512	ILE	N-CA-C	6.02	118.98	112.96
1	B	535	MET	N-CA-C	-5.95	106.18	113.50
1	B	483	MET	CA-C-N	5.90	125.86	119.78
1	B	483	MET	C-N-CA	5.90	125.86	119.78
1	A	535	MET	N-CA-C	-5.89	106.25	113.50
1	A	413	LYS	N-CA-C	5.88	117.69	111.28
1	A	483	MET	CA-C-N	5.88	125.84	119.78
1	A	483	MET	C-N-CA	5.88	125.84	119.78
1	B	334	HIS	N-CA-C	5.86	119.69	112.54
1	A	317	SER	N-CA-C	5.86	118.14	111.11
1	B	512	ILE	N-CA-C	5.80	118.76	112.96
1	A	334	HIS	N-CA-C	5.70	119.49	112.54
1	B	96	LEU	N-CA-C	5.65	118.98	111.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	413	LYS	N-CA-C	5.63	117.41	111.28
1	A	96	LEU	N-CA-C	5.60	118.92	111.75
1	B	317	SER	N-CA-C	5.58	117.81	111.11
1	B	339	GLY	N-CA-C	5.50	119.33	112.73
1	B	549	THR	N-CA-C	5.47	117.29	109.14
1	A	339	GLY	N-CA-C	5.42	119.23	112.73
1	B	563	GLY	CA-C-N	5.28	125.26	120.03
1	B	563	GLY	C-N-CA	5.28	125.26	120.03
1	A	549	THR	N-CA-C	5.27	117.00	109.14
1	A	42	MET	N-CA-C	5.12	119.61	112.90
1	B	461	ASN	N-CA-C	5.12	117.83	109.59
1	B	42	MET	N-CA-C	5.09	119.56	112.90
1	A	461	ASN	N-CA-C	5.05	117.72	109.59
1	A	585	ILE	N-CA-C	5.01	115.13	108.11

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	4667	0	4622	66	0
1	B	4667	0	4621	67	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	149	0	0	3	0
3	B	196	0	0	2	0
All	All	9681	0	9243	133	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 7.

All (133) 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:352:HIS:HA	1:A:355:THR:HG23	1.47	0.97
1:B:352:HIS:HA	1:B:355:THR:HG23	1.47	0.96
1:A:355:THR:C	1:A:356:SER:CA	2.44	0.91
1:B:355:THR:C	1:B:356:SER:CA	2.44	0.89
1:A:551:GLN:HE21	1:A:576:PRO:HG2	1.43	0.83
1:B:551:GLN:HE21	1:B:576:PRO:HG2	1.44	0.82
1:A:367:SER:CA	1:A:368:GLU:N	2.42	0.82
1:B:367:SER:CA	1:B:368:GLU:N	2.42	0.82
1:B:313:TRP:H	1:B:319:GLN:NE2	1.78	0.82
1:A:313:TRP:H	1:A:319:GLN:NE2	1.80	0.78
1:A:489:LEU:HD21	1:A:617:ARG:HD3	1.68	0.76
1:A:211:ILE:O	1:A:212:ASN:HB2	1.86	0.75
1:B:211:ILE:O	1:B:212:ASN:HB2	1.86	0.75
1:B:489:LEU:HD21	1:B:617:ARG:HD3	1.69	0.73
1:B:190:MET:HB3	1:B:191:PRO:HD2	1.71	0.73
1:A:190:MET:HB3	1:A:191:PRO:HD2	1.70	0.73
1:B:551:GLN:NE2	1:B:576:PRO:HG2	2.05	0.72
1:A:319:GLN:HB2	1:A:331:VAL:HG13	1.72	0.71
1:A:551:GLN:NE2	1:A:576:PRO:HG2	2.05	0.71
1:B:319:GLN:HB2	1:B:331:VAL:HG13	1.73	0.68
1:B:313:TRP:H	1:B:319:GLN:HE22	1.42	0.67
1:A:551:GLN:HG2	1:A:553:PRO:HD3	1.80	0.63
1:B:537:GLU:HG3	1:B:541:MET:HE2	1.79	0.63
1:B:50:GLN:HE22	1:B:541:MET:HE1	1.64	0.63
1:A:50:GLN:HE22	1:A:541:MET:HE1	1.63	0.62
1:A:537:GLU:HG3	1:A:541:MET:HE2	1.78	0.62
1:A:313:TRP:H	1:A:319:GLN:HE22	1.45	0.62
1:B:551:GLN:HG2	1:B:553:PRO:HD3	1.80	0.62
1:A:292:ASP:OD2	1:A:295:ARG:NH2	2.34	0.60
1:B:492:LEU:C	1:B:492:LEU:HD23	2.26	0.60
1:A:492:LEU:C	1:A:492:LEU:HD23	2.26	0.60
1:A:211:ILE:HD12	1:A:369:LEU:HD22	1.84	0.59
1:B:211:ILE:HD12	1:B:369:LEU:HD22	1.85	0.59
1:A:153:GLY:HA3	1:A:156:SER:O	2.04	0.58
1:A:415:PHE:CZ	1:A:611:ARG:HG3	2.38	0.58
1:B:292:ASP:OD2	1:B:295:ARG:NH2	2.36	0.58
1:A:151:ALA:HB2	1:A:160:LEU:HD21	1.85	0.58
1:B:153:GLY:HA3	1:B:156:SER:O	2.05	0.57
1:A:479:HIS:HB2	1:A:482:ALA:HB3	1.88	0.55
1:B:415:PHE:CZ	1:B:611:ARG:HG3	2.41	0.55
1:B:479:HIS:HB2	1:B:482:ALA:HB3	1.88	0.55
1:B:479:HIS:ND1	1:B:480:LYS:N	2.55	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:ARG:HD3	1:A:115:GLN:O	2.07	0.55
1:B:479:HIS:CG	1:B:480:LYS:N	2.75	0.55
1:B:114:ARG:HD3	1:B:115:GLN:O	2.07	0.54
1:B:151:ALA:HB2	1:B:160:LEU:HD21	1.89	0.53
1:A:479:HIS:CG	1:A:480:LYS:N	2.76	0.53
1:A:556:MET:HB2	3:A:1098:HOH:O	2.07	0.53
1:A:483:MET:HE3	1:A:487:GLU:HB2	1.90	0.53
1:A:483:MET:HE3	1:A:487:GLU:CB	2.39	0.53
1:B:479:HIS:CD2	1:B:481:ALA:H	2.26	0.53
1:A:479:HIS:CD2	1:A:481:ALA:H	2.27	0.53
1:A:479:HIS:ND1	1:A:480:LYS:N	2.57	0.53
1:B:551:GLN:HE21	1:B:576:PRO:CG	2.20	0.53
1:B:483:MET:HE3	1:B:487:GLU:CB	2.39	0.53
1:B:599:PHE:O	1:B:603:VAL:HG23	2.09	0.53
1:B:483:MET:HE3	1:B:487:GLU:HB2	1.91	0.52
1:A:599:PHE:O	1:A:603:VAL:HG23	2.10	0.52
1:A:111:ILE:HD12	1:A:111:ILE:N	2.24	0.52
1:B:352:HIS:HA	1:B:355:THR:CG2	2.31	0.52
1:B:111:ILE:HD12	1:B:111:ILE:N	2.25	0.51
1:A:177:HIS:CD2	1:A:178:THR:HG23	2.45	0.51
1:B:177:HIS:CD2	1:B:178:THR:HG23	2.46	0.51
1:A:352:HIS:HA	1:A:355:THR:CG2	2.31	0.51
1:A:400:ASN:HD21	1:A:590:SER:H	1.59	0.49
1:A:419:GLN:HG3	1:A:617:ARG:NH2	2.28	0.48
1:B:441:ARG:HG3	1:B:441:ARG:HH11	1.79	0.48
1:B:419:GLN:HG3	1:B:617:ARG:NH2	2.29	0.48
1:A:320:PHE:CD2	1:A:330:VAL:HG22	2.48	0.48
1:B:480:LYS:O	1:B:481:ALA:HB3	2.13	0.48
1:A:480:LYS:O	1:A:481:ALA:HB3	2.14	0.47
1:A:57:ILE:HG23	1:A:532:GLU:HG2	1.97	0.47
1:B:400:ASN:HD21	1:B:590:SER:H	1.61	0.47
1:A:441:ARG:HH11	1:A:441:ARG:HG3	1.80	0.47
1:A:99:ARG:N	1:A:99:ARG:HD2	2.30	0.46
1:B:57:ILE:HG23	1:B:532:GLU:HG2	1.98	0.46
1:A:374:ARG:NH2	3:A:1030:HOH:O	2.43	0.46
1:B:114:ARG:HH11	1:B:114:ARG:HG3	1.80	0.46
1:B:99:ARG:HD2	1:B:99:ARG:N	2.29	0.46
1:A:425:GLY:O	1:A:429:VAL:HG23	2.16	0.46
1:B:320:PHE:CD2	1:B:330:VAL:HG22	2.49	0.46
1:A:114:ARG:HH11	1:A:114:ARG:HG3	1.81	0.45
1:A:551:GLN:HE21	1:A:576:PRO:CG	2.19	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:HIS:HD2	1:A:178:THR:HG23	1.82	0.44
1:B:177:HIS:HD2	1:B:178:THR:HG23	1.82	0.44
1:B:114:ARG:HG3	1:B:114:ARG:NH1	2.33	0.43
1:A:464:SER:HA	1:A:499:GLN:HE22	1.83	0.43
1:B:425:GLY:O	1:B:429:VAL:HG23	2.17	0.43
1:B:537:GLU:HG3	1:B:541:MET:CE	2.46	0.43
1:A:464:SER:HA	1:A:499:GLN:NE2	2.33	0.43
1:B:270:LEU:HD22	3:B:1068:HOH:O	2.18	0.43
1:B:464:SER:HA	1:B:499:GLN:NE2	2.34	0.43
1:B:212:ASN:HB3	1:B:213:PHE:H	1.60	0.43
1:A:470:LEU:O	1:A:474:GLN:HG3	2.19	0.43
1:B:374:ARG:NH2	3:B:1047:HOH:O	2.45	0.43
1:A:114:ARG:HG3	1:A:114:ARG:NH1	2.34	0.43
1:A:445:THR:HG23	1:A:547:LEU:HB3	2.01	0.43
1:B:530:PRO:HA	1:B:531:PRO:HD3	1.89	0.42
1:A:106:SER:O	1:A:108:PRO:HD3	2.18	0.42
1:B:153:GLY:CA	1:B:156:SER:O	2.67	0.42
1:B:547:LEU:HD22	1:B:549:THR:CG2	2.49	0.42
1:A:212:ASN:HB3	1:A:213:PHE:H	1.61	0.42
1:A:421:TYR:HB3	1:A:496:MET:HE2	2.01	0.42
1:B:483:MET:HA	1:B:484:PRO:HD3	1.88	0.42
1:B:410:ASN:O	1:B:411:TYR:HB3	2.20	0.42
1:A:20:GLU:HB3	1:A:23:LEU:HD12	2.00	0.42
1:A:340:ILE:O	1:A:344:GLN:HG2	2.19	0.42
1:A:547:LEU:HD22	1:A:549:THR:CG2	2.48	0.42
1:B:20:GLU:HB3	1:B:23:LEU:HD12	2.01	0.42
1:B:214:ARG:NH2	1:B:369:LEU:HD21	2.34	0.42
1:A:99:ARG:NH2	1:A:176:GLY:O	2.53	0.41
1:A:102:LEU:N	1:A:103:PRO:CD	2.83	0.41
1:A:238:ARG:NH2	3:A:1092:HOH:O	2.51	0.41
1:A:500:THR:O	1:A:504:VAL:HG23	2.20	0.41
1:B:99:ARG:NH2	1:B:176:GLY:O	2.53	0.41
1:B:464:SER:HA	1:B:499:GLN:HE22	1.84	0.41
1:A:153:GLY:CA	1:A:156:SER:O	2.68	0.41
1:A:214:ARG:NH2	1:A:369:LEU:HD21	2.35	0.41
1:A:130:LEU:C	1:A:130:LEU:HD23	2.45	0.41
1:B:26:LEU:HA	1:B:27:PRO:HD3	1.90	0.41
1:B:421:TYR:HB3	1:B:496:MET:HE2	2.01	0.41
1:B:340:ILE:O	1:B:344:GLN:HG2	2.20	0.41
1:B:106:SER:O	1:B:108:PRO:HD3	2.21	0.41
1:A:26:LEU:HA	1:A:27:PRO:HD3	1.90	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:50:GLN:OE1	1:B:541:MET:HE3	2.21	0.41
1:A:321:VAL:O	1:A:328:CYS:HA	2.21	0.41
1:B:194:GLU:O	1:B:209:VAL:HG22	2.21	0.41
1:B:214:ARG:HH22	1:B:369:LEU:HD21	1.86	0.41
1:B:470:LEU:O	1:B:474:GLN:HG3	2.21	0.41
1:B:479:HIS:CD2	1:B:481:ALA:N	2.89	0.41
1:A:396:ARG:O	1:A:400:ASN:HB2	2.20	0.40
1:B:445:THR:HG23	1:B:547:LEU:HB3	2.02	0.40
1:A:537:GLU:HG3	1:A:541:MET:CE	2.47	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	584/644 (91%)	559 (96%)	24 (4%)	1 (0%)	43	63
1	B	584/644 (91%)	558 (96%)	25 (4%)	1 (0%)	43	63
All	All	1168/1288 (91%)	1117 (96%)	49 (4%)	2 (0%)	43	63

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	212	ASN
1	B	212	ASN

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	516/565 (91%)	504 (98%)	12 (2%)	44	72
1	B	516/565 (91%)	504 (98%)	12 (2%)	44	72
All	All	1032/1130 (91%)	1008 (98%)	24 (2%)	44	72

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	GLU
1	A	100	LEU
1	A	150	TRP
1	A	212	ASN
1	A	244	LEU
1	A	260	LEU
1	A	279	LEU
1	A	331	VAL
1	A	400	ASN
1	A	407	LYS
1	A	419	GLN
1	A	547	LEU
1	B	20	GLU
1	B	100	LEU
1	B	150	TRP
1	B	212	ASN
1	B	244	LEU
1	B	260	LEU
1	B	279	LEU
1	B	331	VAL
1	B	400	ASN
1	B	407	LYS
1	B	419	GLN
1	B	547	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	121	ASN
1	A	177	HIS

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Mol	Chain	Res	Type
1	A	179	GLN
1	A	185	GLN
1	A	202	ASN
1	A	234	ASN
1	A	298	GLN
1	A	319	GLN
1	A	384	GLN
1	A	395	GLN
1	A	400	ASN
1	A	410	ASN
1	A	455	GLN
1	A	474	GLN
1	A	493	GLN
1	A	499	GLN
1	A	551	GLN
1	B	33	GLN
1	B	121	ASN
1	B	177	HIS
1	B	202	ASN
1	B	298	GLN
1	B	319	GLN
1	B	384	GLN
1	B	400	ASN
1	B	410	ASN
1	B	455	GLN
1	B	474	GLN
1	B	493	GLN
1	B	499	GLN
1	B	551	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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

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	600/644 (93%)	0.81	43 (7%)	21 19	12, 29, 73, 110	0
1	B	600/644 (93%)	0.62	38 (6%)	26 23	11, 24, 66, 108	0
All	All	1200/1288 (93%)	0.71	81 (6%)	23 20	11, 26, 71, 110	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	361	VAL	36.9
1	B	366	VAL	35.0
1	B	364	ASP	34.2
1	B	365	SER	33.9
1	A	364	ASP	32.9
1	A	365	SER	32.8
1	A	366	VAL	32.0
1	A	357	ASN	31.4
1	B	357	ASN	29.2
1	A	362	ARG	28.7
1	B	361	VAL	27.9
1	B	363	ALA	27.6
1	B	359	LYS	26.3
1	A	363	ALA	26.2
1	B	362	ARG	26.2
1	A	360	LEU	26.0
1	A	359	LYS	25.5
1	A	356	SER	25.3
1	A	358	LYS	24.2
1	B	360	LEU	23.7
1	B	367	SER	23.2
1	B	356	SER	22.8
1	A	367	SER	22.5
1	B	358	LYS	22.4

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Mol	Chain	Res	Type	RSRZ
1	A	189	ILE	7.4
1	B	189	ILE	7.0
1	A	191	PRO	5.5
1	B	191	PRO	5.5
1	B	213	PHE	5.2
1	A	213	PHE	5.1
1	B	187	SER	4.7
1	A	190	MET	4.7
1	B	211	ILE	4.6
1	A	211	ILE	4.5
1	B	190	MET	4.4
1	B	355	THR	4.0
1	A	188	SER	3.8
1	A	482	ALA	3.7
1	A	18	TRP	3.7
1	A	355	THR	3.6
1	A	187	SER	3.5
1	B	18	TRP	3.5
1	B	482	ALA	3.5
1	A	485	ALA	3.4
1	B	214	ARG	3.4
1	B	617	ARG	3.4
1	B	188	SER	3.4
1	A	193	PRO	3.4
1	B	480	LYS	3.3
1	B	354	MET	3.3
1	B	616	SER	3.2
1	A	479	HIS	3.1
1	B	186	LYS	3.1
1	A	617	ARG	3.0
1	A	481	ALA	3.0
1	B	481	ALA	3.0
1	A	480	LYS	3.0
1	B	479	HIS	3.0
1	B	193	PRO	2.7
1	A	354	MET	2.6
1	A	486	SER	2.6
1	A	186	LYS	2.6
1	B	177	HIS	2.5
1	A	477	THR	2.5
1	A	410	ASN	2.4
1	A	192	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	212	ASN	2.4
1	A	352	HIS	2.3
1	B	192	GLU	2.3
1	A	613	LEU	2.3
1	A	484	PRO	2.2
1	B	615	SER	2.2
1	B	210	VAL	2.2
1	A	483	MET	2.1
1	B	612	ASP	2.1
1	A	490	GLN	2.1
1	A	529	GLU	2.1
1	A	493	GLN	2.1
1	B	212	ASN	2.1
1	B	52	ARG	2.1
1	A	19	GLU	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NA	A	1002	1/1	0.94	0.06	24,24,24,24	0
2	NA	B	1001	1/1	0.94	0.05	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.