



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 2NQ5 / pdb_00002nq5
Title : Crystal structure of methyltransferase from Streptococcus mutans
Authors : Fedorov, A.A.; Fedorov, E.V.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2006-10-30
Resolution : 1.90 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

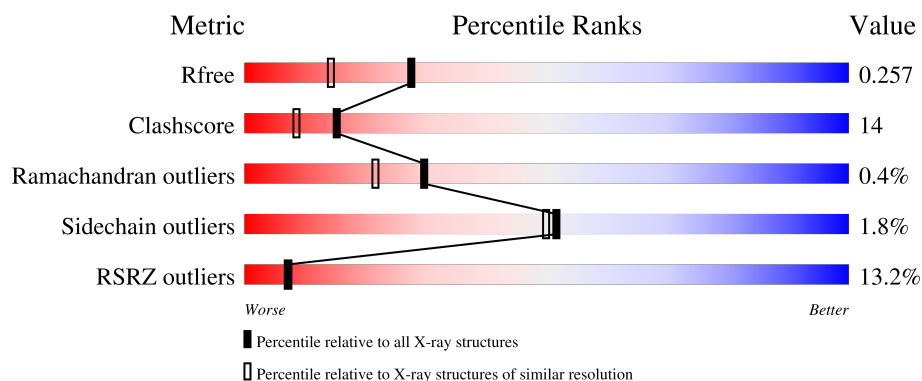
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	755	12% 68% 24% 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5980 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 5-methyltetrahydropteroyltriglutamate--homocysteine methyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	706	Total	C	N	O	S	0	0	0
			5632	3623	943	1059	7			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	cloning artifact	UNP Q8CWX6
A	0	SER	-	cloning artifact	UNP Q8CWX6
A	1	LEU	-	cloning artifact	UNP Q8CWX6
A	746	GLU	-	cloning artifact	UNP Q8CWX6
A	747	GLY	-	cloning artifact	UNP Q8CWX6
A	748	HIS	-	expression tag	UNP Q8CWX6
A	749	HIS	-	expression tag	UNP Q8CWX6
A	750	HIS	-	expression tag	UNP Q8CWX6
A	751	HIS	-	expression tag	UNP Q8CWX6
A	752	HIS	-	expression tag	UNP Q8CWX6
A	753	HIS	-	expression tag	UNP Q8CWX6

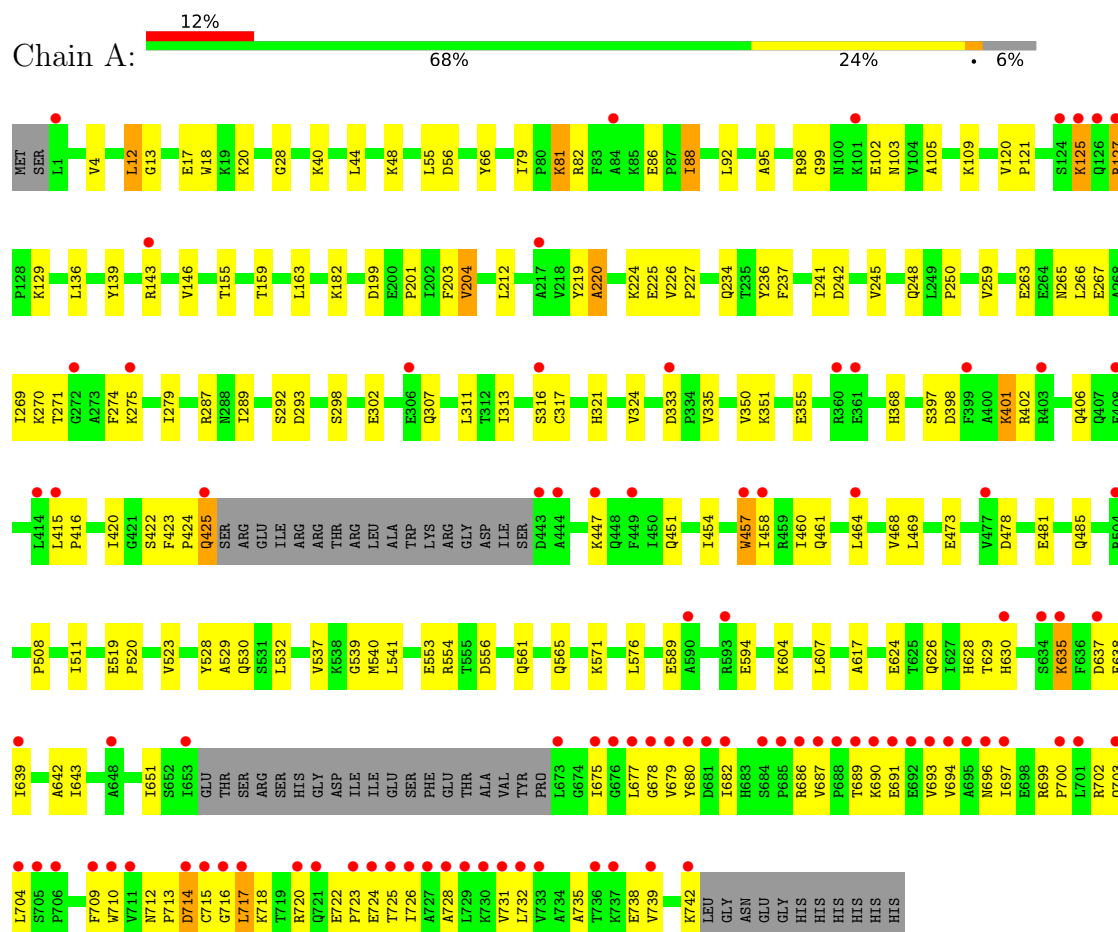
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	348	Total	O	0	0
			348	348		

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: 5-methyltetrahydropteroyltriglutamate--homocysteine methyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.29Å 99.19Å 78.62Å 90.00° 109.25° 90.00°	Depositor
Resolution (Å)	24.96 – 1.90 24.96 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.5 (24.96-1.90) 97.5 (24.96-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.62 (at 1.88Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.228 , 0.258 0.228 , 0.257	Depositor DCC
R_{free} test set	3028 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	18.2	Xtriage
Anisotropy	0.268	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 44.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.029 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5980	wwPDB-VP
Average B, all atoms (Å ²)	24.0	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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	0.38	0/5752	0.89	17/7801 (0.2%)

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	1

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	324	VAL	N-CA-C	9.57	119.53	110.53
1	A	204	VAL	N-CA-C	-8.59	104.77	113.10
1	A	12	LEU	N-CA-C	7.10	119.92	111.33
1	A	13	GLY	N-CA-C	-6.06	105.04	112.68
1	A	234	GLN	N-CA-C	6.04	118.50	109.25
1	A	146	VAL	N-CA-C	6.01	118.67	112.90
1	A	220	ALA	N-CA-C	-5.85	104.98	111.36
1	A	226	VAL	CA-C-N	5.70	125.82	119.32
1	A	226	VAL	C-N-CA	5.70	125.82	119.32
1	A	539	GLY	N-CA-C	-5.64	102.44	112.54
1	A	508	PRO	CA-C-N	-5.57	115.00	120.52
1	A	508	PRO	C-N-CA	-5.57	115.00	120.52
1	A	523	VAL	N-CA-C	5.53	115.73	110.42
1	A	289	ILE	N-CA-C	5.26	120.28	109.34
1	A	321	HIS	N-CA-C	5.18	119.23	113.02
1	A	293	ASP	N-CA-C	-5.11	99.00	107.99
1	A	469	LEU	N-CA-C	5.01	117.87	110.46

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	219	TYR	Sidechain

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	5632	0	5660	153	0
2	A	348	0	0	3	0
All	All	5980	0	5660	153	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 14.

All (153) 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:143:ARG:HH11	1:A:143:ARG:HG3	1.36	0.91
1:A:4:VAL:HG23	1:A:55:LEU:HA	1.52	0.91
1:A:125:LYS:H	1:A:125:LYS:HD2	1.39	0.88
1:A:553:GLU:HG2	2:A:978:HOH:O	1.75	0.87
1:A:630:HIS:NE2	1:A:715:CYS:SG	2.53	0.81
1:A:717:LEU:HD23	1:A:717:LEU:O	1.83	0.78
1:A:425:GLN:HE21	1:A:425:GLN:N	1.82	0.77
1:A:275:LYS:O	1:A:275:LYS:HD3	1.83	0.77
1:A:720:ARG:HH12	1:A:724:GLU:HB3	1.48	0.77
1:A:313:ILE:HG22	1:A:350:VAL:HG13	1.67	0.76
1:A:717:LEU:HD21	1:A:725:THR:HG21	1.68	0.75
1:A:458:ILE:HD12	1:A:532:LEU:HD12	1.70	0.74
1:A:626:GLN:HE21	1:A:628:HIS:HE1	1.35	0.73
1:A:722:GLU:HB3	1:A:723:PRO:HD3	1.71	0.72
1:A:626:GLN:NE2	1:A:628:HIS:HE1	1.89	0.71
1:A:127:ARG:O	1:A:127:ARG:HD2	1.91	0.71
1:A:511:ILE:O	1:A:554:ARG:HG3	1.92	0.70
1:A:626:GLN:HE21	1:A:628:HIS:CE1	2.10	0.69
1:A:637:ASP:OD1	1:A:638:GLU:HG2	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:481:GLU:O	1:A:485:GLN:HG3	1.94	0.67
1:A:397:SER:OG	1:A:401:LYS:HE2	1.95	0.67
1:A:292:SER:O	1:A:368:HIS:HE1	1.78	0.66
1:A:457:TRP:O	1:A:461:GLN:HG3	1.95	0.66
1:A:127:ARG:HE	1:A:129:LYS:HD3	1.60	0.66
1:A:245:VAL:O	1:A:248:GLN:HG2	1.96	0.66
1:A:635:LYS:HB2	1:A:635:LYS:NZ	2.11	0.65
1:A:143:ARG:HH11	1:A:143:ARG:CG	2.10	0.65
1:A:88:ILE:HD12	1:A:88:ILE:N	2.11	0.65
1:A:447:LYS:O	1:A:451:GLN:HG3	1.96	0.65
1:A:263:GLU:O	1:A:267:GLU:HG3	1.98	0.64
1:A:159:THR:O	1:A:163:LEU:HD13	1.98	0.63
1:A:473:GLU:OE2	1:A:540:MET:HG2	1.99	0.63
1:A:4:VAL:CG2	1:A:55:LEU:HA	2.28	0.62
1:A:639:ILE:O	1:A:643:ILE:HG12	1.99	0.62
1:A:717:LEU:HD21	1:A:725:THR:CG2	2.29	0.62
1:A:738:GLU:O	1:A:742:LYS:HD3	2.00	0.61
1:A:333:ASP:OD1	1:A:335:VAL:HG12	1.99	0.61
1:A:607:LEU:HD13	1:A:639:ILE:HG23	1.82	0.60
1:A:28:GLY:HA2	2:A:1090:HOH:O	2.01	0.59
1:A:125:LYS:HD2	1:A:125:LYS:N	2.16	0.59
1:A:604:LYS:HG2	2:A:1009:HOH:O	2.04	0.58
1:A:82:ARG:HB2	1:A:103:ASN:ND2	2.18	0.58
1:A:425:GLN:OE1	1:A:718:LYS:HE3	2.04	0.58
1:A:699:ARG:O	1:A:702:ARG:HB3	2.02	0.58
1:A:298:SER:O	1:A:302:GLU:HG3	2.04	0.57
1:A:136:LEU:C	1:A:136:LEU:HD13	2.29	0.57
1:A:690:LYS:HG2	1:A:731:VAL:HG23	1.85	0.57
1:A:626:GLN:HG2	1:A:628:HIS:CE1	2.40	0.56
1:A:351:LYS:O	1:A:355:GLU:HG3	2.05	0.56
1:A:682:ILE:HD12	1:A:682:ILE:C	2.31	0.56
1:A:98:ARG:C	1:A:105:ALA:HB2	2.31	0.56
1:A:424:PRO:HD3	1:A:717:LEU:CD2	2.36	0.56
1:A:81:LYS:HE3	1:A:81:LYS:HA	1.87	0.55
1:A:143:ARG:HG3	1:A:143:ARG:NH1	2.14	0.55
1:A:398:ASP:O	1:A:402:ARG:HG3	2.06	0.55
1:A:704:LEU:HD12	1:A:704:LEU:N	2.21	0.55
1:A:201:PRO:O	1:A:204:VAL:HG22	2.06	0.55
1:A:275:LYS:HD3	1:A:275:LYS:C	2.31	0.55
1:A:17:GLU:HA	1:A:20:LYS:HE3	1.87	0.55
1:A:267:GLU:O	1:A:271:THR:HG23	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:720:ARG:HH12	1:A:724:GLU:CB	2.20	0.54
1:A:109:LYS:HG3	1:A:120:VAL:HG21	1.90	0.53
1:A:693:VAL:O	1:A:697:ILE:HG22	2.08	0.53
1:A:720:ARG:HH11	1:A:720:ARG:HG3	1.73	0.53
1:A:678:GLY:HA2	1:A:712:ASN:O	2.09	0.52
1:A:4:VAL:HG11	1:A:351:LYS:HA	1.92	0.52
1:A:127:ARG:HH11	1:A:127:ARG:HG2	1.73	0.52
1:A:460:ILE:O	1:A:464:LEU:HG	2.10	0.51
1:A:287:ARG:HG3	1:A:317:CYS:SG	2.51	0.51
1:A:415:LEU:HD13	1:A:697:ILE:HD11	1.92	0.51
1:A:630:HIS:HE2	1:A:715:CYS:HA	1.76	0.51
1:A:677:LEU:HD11	1:A:709:PHE:HD2	1.76	0.51
1:A:423:PHE:HA	1:A:717:LEU:HD22	1.93	0.51
1:A:425:GLN:HE22	1:A:718:LYS:HG2	1.76	0.51
1:A:397:SER:CB	1:A:401:LYS:HE2	2.42	0.50
1:A:704:LEU:HD12	1:A:704:LEU:H	1.77	0.49
1:A:313:ILE:CG2	1:A:350:VAL:HG13	2.40	0.49
1:A:422:SER:O	1:A:717:LEU:HB3	2.12	0.49
1:A:702:ARG:HG3	1:A:703:GLN:HG3	1.94	0.49
1:A:629:THR:O	1:A:651:ILE:HA	2.13	0.48
1:A:561:GLN:O	1:A:565:GLN:HG3	2.13	0.48
1:A:677:LEU:HD11	1:A:709:PHE:CD2	2.47	0.48
1:A:139:TYR:CZ	1:A:143:ARG:HD3	2.49	0.48
1:A:451:GLN:HB3	1:A:528:TYR:CE2	2.48	0.47
1:A:680:TYR:HD1	1:A:713:PRO:HB3	1.79	0.47
1:A:40:LYS:O	1:A:44:LEU:HG	2.15	0.47
1:A:4:VAL:HG22	1:A:56:ASP:OD2	2.15	0.47
1:A:224:LYS:NZ	1:A:225:GLU:OE2	2.47	0.47
1:A:709:PHE:CD1	1:A:709:PHE:N	2.83	0.47
1:A:266:LEU:HG	1:A:270:LYS:HE2	1.97	0.47
1:A:651:ILE:O	1:A:651:ILE:HG23	2.15	0.47
1:A:478:ASP:HB3	1:A:481:GLU:HB3	1.96	0.47
1:A:182:LYS:C	1:A:182:LYS:HD3	2.40	0.46
1:A:704:LEU:H	1:A:704:LEU:CD1	2.28	0.46
1:A:689:THR:O	1:A:693:VAL:HG22	2.15	0.46
1:A:732:LEU:C	1:A:732:LEU:HD23	2.40	0.46
1:A:220:ALA:HA	1:A:250:PRO:HG3	1.97	0.46
1:A:79:ILE:HD13	1:A:92:LEU:HD21	1.97	0.46
1:A:44:LEU:O	1:A:48:LYS:HG3	2.16	0.46
1:A:88:ILE:N	1:A:88:ILE:CD1	2.79	0.46
1:A:680:TYR:CD1	1:A:713:PRO:HB3	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:735:ALA:O	1:A:739:VAL:HG23	2.16	0.45
1:A:639:ILE:O	1:A:639:ILE:HG22	2.16	0.45
1:A:691:GLU:C	1:A:694:VAL:HG12	2.41	0.45
1:A:259:VAL:HB	1:A:316:SER:OG	2.16	0.45
1:A:720:ARG:NH1	1:A:724:GLU:HB3	2.26	0.45
1:A:236:TYR:O	1:A:237:PHE:HB2	2.15	0.45
1:A:12:LEU:HA	1:A:18:TRP:HB3	1.98	0.45
1:A:265:ASN:O	1:A:269:ILE:HG12	2.17	0.44
1:A:458:ILE:HD11	1:A:529:ALA:HA	1.99	0.44
1:A:571:LYS:HB2	1:A:617:ALA:HA	1.99	0.44
1:A:686:ARG:O	1:A:720:ARG:NH2	2.50	0.44
1:A:675:ILE:HD11	1:A:709:PHE:CE2	2.53	0.44
1:A:691:GLU:O	1:A:694:VAL:HG12	2.18	0.44
1:A:687:VAL:O	1:A:687:VAL:HG13	2.18	0.44
1:A:530:GLN:NE2	1:A:537:VAL:H	2.16	0.44
1:A:554:ARG:HB3	1:A:556:ASP:OD1	2.18	0.44
1:A:86:GLU:HB2	1:A:92:LEU:HD13	2.00	0.44
1:A:127:ARG:HG2	1:A:127:ARG:NH1	2.33	0.44
1:A:127:ARG:HD2	1:A:127:ARG:C	2.43	0.43
1:A:225:GLU:C	1:A:227:PRO:HD3	2.43	0.43
1:A:270:LYS:HA	1:A:307:GLN:HE21	1.83	0.43
1:A:589:GLU:OE1	1:A:589:GLU:HA	2.18	0.43
1:A:540:MET:O	1:A:541:LEU:HD23	2.18	0.43
1:A:415:LEU:N	1:A:416:PRO:HD3	2.31	0.43
1:A:519:GLU:HB2	1:A:520:PRO:HD2	2.00	0.43
1:A:95:ALA:O	1:A:99:GLY:N	2.49	0.43
1:A:722:GLU:O	1:A:726:ILE:HG22	2.18	0.43
1:A:728:ALA:O	1:A:731:VAL:HG12	2.18	0.43
1:A:212:LEU:HD11	1:A:242:ASP:OD2	2.19	0.43
1:A:274:PHE:HB3	1:A:279:ILE:HD11	2.01	0.43
1:A:690:LYS:O	1:A:694:VAL:HG12	2.19	0.43
1:A:406:GLN:HG2	1:A:624:GLU:O	2.18	0.42
1:A:639:ILE:HG22	1:A:642:ALA:HB3	2.00	0.42
1:A:699:ARG:N	1:A:700:PRO:HD2	2.33	0.42
1:A:635:LYS:HB2	1:A:635:LYS:HZ2	1.80	0.42
1:A:702:ARG:HD2	1:A:702:ARG:O	2.18	0.42
1:A:99:GLY:N	1:A:105:ALA:HB2	2.35	0.42
1:A:454:ILE:O	1:A:458:ILE:HG13	2.19	0.42
1:A:203:PHE:O	1:A:241:ILE:HG12	2.19	0.42
1:A:451:GLN:HB3	1:A:528:TYR:CZ	2.55	0.42
1:A:425:GLN:HE21	1:A:425:GLN:H	1.62	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:682:ILE:HG21	1:A:716:GLY:O	2.19	0.41
1:A:594:GLU:CD	1:A:594:GLU:N	2.78	0.41
1:A:155:THR:HA	1:A:199:ASP:HB2	2.01	0.41
1:A:704:LEU:N	1:A:704:LEU:CD1	2.84	0.41
1:A:420:ILE:HG13	1:A:714:ASP:O	2.21	0.41
1:A:679:VAL:HG12	1:A:696:ASN:HB3	2.03	0.41
1:A:630:HIS:CD2	1:A:715:CYS:SG	3.14	0.41
1:A:693:VAL:HG23	1:A:735:ALA:CB	2.51	0.41
1:A:120:VAL:HA	1:A:121:PRO:HD3	1.94	0.40
1:A:143:ARG:CG	1:A:143:ARG:NH1	2.72	0.40
1:A:468:VAL:HG21	1:A:710:TRP:CZ3	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	700/755 (93%)	676 (97%)	21 (3%)	3 (0%)	30	22

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	714	ASP
1	A	717	LEU
1	A	66	TYR

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	608/651 (93%)	597 (98%)	11 (2%)	51	50

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

Mol	Chain	Res	Type
1	A	81	LYS
1	A	88	ILE
1	A	102	GLU
1	A	125	LYS
1	A	127	ARG
1	A	311	LEU
1	A	401	LYS
1	A	425	GLN
1	A	457	TRP
1	A	576	LEU
1	A	635	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	ASN
1	A	50	GLN
1	A	103	ASN
1	A	115	ASN
1	A	234	GLN
1	A	314	GLN
1	A	356	HIS
1	A	368	HIS
1	A	425	GLN
1	A	530	GLN
1	A	586	GLN
1	A	628	HIS
1	A	696	ASN
1	A	703	GLN

5.3.3 RNA ⓘ

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 ⓘ

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	706/755 (93%)	0.64	93 (13%) 7 7	12, 21, 43, 52	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	715	CYS	6.4
1	A	673	LEU	6.3
1	A	697	ILE	6.2
1	A	716	GLY	6.2
1	A	1	LEU	5.9
1	A	675	ILE	5.9
1	A	679	VAL	5.6
1	A	692	GLU	5.3
1	A	703	GLN	5.2
1	A	449	PHE	4.8
1	A	680	TYR	4.7
1	A	704	LEU	4.7
1	A	720	ARG	4.6
1	A	726	ILE	4.4
1	A	639	ILE	4.4
1	A	711	VAL	4.4
1	A	414	LEU	4.3
1	A	635	LYS	4.3
1	A	723	PRO	4.3
1	A	695	ALA	4.2
1	A	694	VAL	4.1
1	A	457	TRP	4.0
1	A	634	SER	4.0
1	A	739	VAL	3.9
1	A	126	GLN	3.9
1	A	717	LEU	3.8
1	A	678	GLY	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	689	THR	3.7
1	A	721	GLN	3.6
1	A	700	PRO	3.6
1	A	685	PRO	3.4
1	A	714	ASP	3.2
1	A	732	LEU	3.2
1	A	653	ILE	3.1
1	A	705	SER	3.1
1	A	677	LEU	3.1
1	A	693	VAL	3.1
1	A	731	VAL	3.0
1	A	701	LEU	3.0
1	A	684	SER	3.0
1	A	399	PHE	2.9
1	A	590	ALA	2.9
1	A	443	ASP	2.9
1	A	696	ASN	2.8
1	A	630	HIS	2.8
1	A	729	LEU	2.8
1	A	124	SER	2.7
1	A	686	ARG	2.7
1	A	408	GLU	2.7
1	A	709	PHE	2.7
1	A	706	PRO	2.7
1	A	690	LYS	2.6
1	A	415	LEU	2.6
1	A	447	LYS	2.6
1	A	127	ARG	2.5
1	A	125	LYS	2.5
1	A	730	LYS	2.5
1	A	687	VAL	2.5
1	A	333	ASP	2.5
1	A	403	ARG	2.5
1	A	504	ARG	2.5
1	A	691	GLU	2.5
1	A	444	ALA	2.5
1	A	682	ILE	2.4
1	A	217	ALA	2.4
1	A	464	LEU	2.3
1	A	458	ILE	2.3
1	A	688	PRO	2.3
1	A	681	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	648	ALA	2.3
1	A	727	ALA	2.3
1	A	733	VAL	2.3
1	A	361	GLU	2.3
1	A	724	GLU	2.3
1	A	676	GLY	2.2
1	A	316	SER	2.2
1	A	425	GLN	2.2
1	A	736	THR	2.2
1	A	101	LYS	2.2
1	A	710	TRP	2.2
1	A	477	VAL	2.2
1	A	275	LYS	2.2
1	A	360	ARG	2.2
1	A	737	LYS	2.2
1	A	725	THR	2.1
1	A	143	ARG	2.1
1	A	742	LYS	2.1
1	A	84	ALA	2.1
1	A	637	ASP	2.1
1	A	272	GLY	2.1
1	A	728	ALA	2.1
1	A	306	GLU	2.0
1	A	593	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.