



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 12:54 pm GMT

PDB ID : 2B7L
Title : Crystal Structure of CTP:Glycerol-3-Phosphate Cytidylyltransferase from Staphylococcus aureus
Authors : Fong, D.H.; Yim, V.C.-N.; D'Elia, M.A.; Brown, E.D.; Berghuis, A.M.
Deposited on : 2005-10-04
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

<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

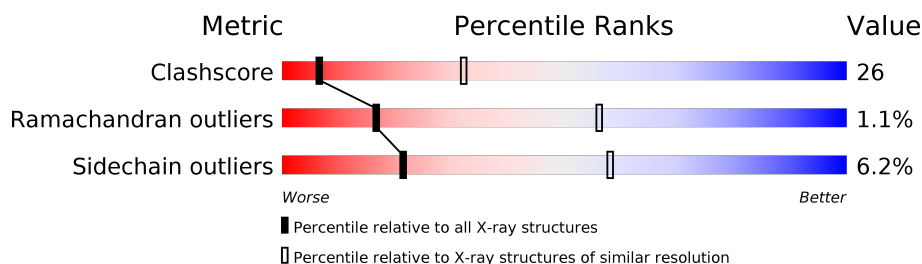
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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))
Clashscore	112137	2037 (3.00-3.00)
Ramachandran outliers	110173	1973 (3.00-3.00)
Sidechain outliers	110143	1976 (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 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	132	
1	B	132	
1	C	132	
1	D	132	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3880 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 glycerol-3-phosphate cytidylyltransferase.

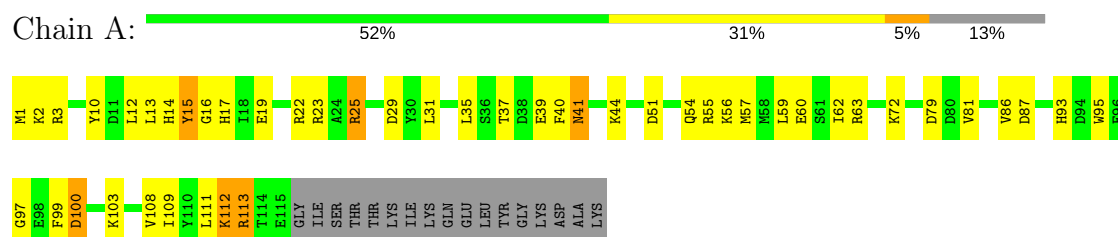
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	115	Total	C	N	O	S	9	0	0
			982	632	161	183	6			
1	B	113	Total	C	N	O	S	11	0	0
			966	623	159	178	6			
1	C	113	Total	C	N	O	S	10	0	0
			966	623	159	178	6			
1	D	113	Total	C	N	O	S	8	0	0
			966	623	159	178	6			

3 Residue-property plots

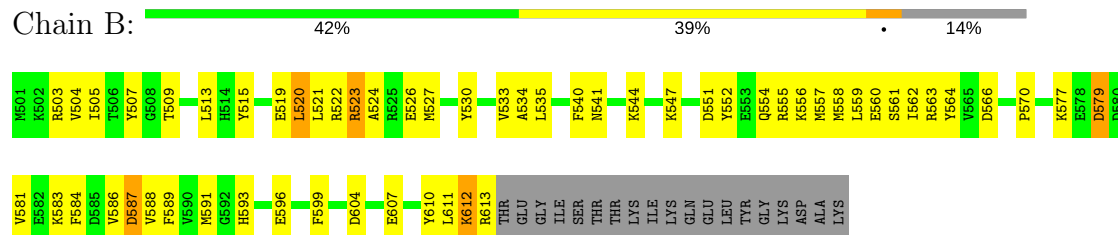
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

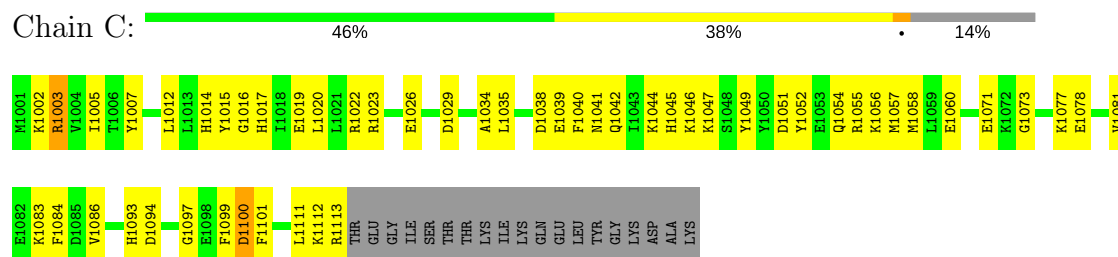
• Molecule 1: glycerol-3-phosphate cytidylyltransferase



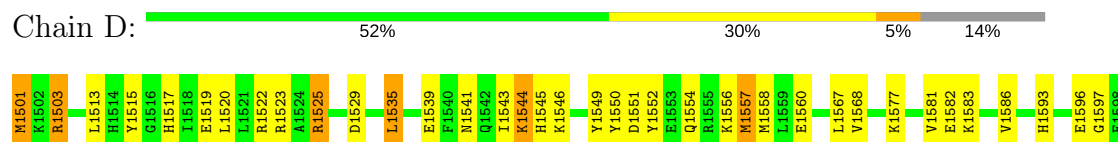
• Molecule 1: glycerol-3-phosphate cytidylyltransferase



• Molecule 1: glycerol-3-phosphate cytidylyltransferase



• Molecule 1: glycerol-3-phosphate cytidylyltransferase



F1599	D1600	F1601	L1602	K1603	D1604	Y1610	L1611	K1612	R1613	THR	GLU	GLY	ILE	SER	THR	THR	LYS	ILE	LYS	GLN	GLU	LEU	TYR	GLY	LYS	ASP	ALA	LYS
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4 Data and refinement statistics

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

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	92.15Å 92.15Å 156.12Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.08 – 3.00	Depositor
% Data completeness (in resolution range)	92.4 (46.08-3.00)	Depositor
R_{merge}	0.14	Depositor
R_{sym}	0.14	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.239 , 0.275	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3880	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

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	0.42	0/1004	0.63	0/1347
1	B	0.43	0/988	0.62	0/1325
1	C	0.44	0/988	0.64	0/1325
1	D	0.41	0/988	0.64	0/1325
All	All	0.43	0/3968	0.63	0/5322

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	982	0	959	61	0
1	B	966	0	943	48	0
1	C	966	0	943	70	0
1	D	966	0	943	40	0
All	All	3880	0	3788	200	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 26.

All (200) 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:112:LYS:H	1:A:112:LYS:HD3	1.04	1.09
1:D:1520:LEU:HD22	1:D:1611:LEU:HD12	1.32	1.06
1:C:1023:ARG:HH22	1:C:1112:LYS:HG3	1.24	1.02
1:A:23:ARG:HH11	1:A:112:LYS:HE2	1.17	1.02
1:D:1525:ARG:HH21	1:D:1529:ASP:HA	1.28	0.96
1:A:112:LYS:HG2	1:A:113:ARG:H	1.35	0.91
1:C:1023:ARG:HH12	1:C:1112:LYS:HE3	1.34	0.90
1:C:1014:HIS:CE1	1:C:1016:GLY:HA3	2.08	0.88
1:B:523:ARG:HG3	1:B:523:ARG:HH11	1.38	0.88
1:B:579:ASP:O	1:B:583:LYS:HG2	1.76	0.85
1:A:112:LYS:N	1:A:112:LYS:HD3	1.87	0.85
1:C:1038:ASP:O	1:C:1041:ASN:HB2	1.81	0.80
1:A:23:ARG:NH1	1:A:112:LYS:HE2	1.97	0.79
1:B:519:GLU:OE2	1:B:523:ARG:HD3	1.84	0.78
1:C:1112:LYS:HG3	1:C:1113:ARG:H	1.49	0.77
1:D:1556:LYS:O	1:D:1560:GLU:HG3	1.84	0.76
1:B:524:ALA:HA	1:B:527:MET:HE3	1.67	0.76
1:B:556:LYS:O	1:B:560:GLU:HG3	1.86	0.76
1:A:56:LYS:O	1:A:60:GLU:HG3	1.85	0.76
1:C:1023:ARG:NH1	1:C:1112:LYS:HE3	2.02	0.75
1:A:3:ARG:HH11	1:A:3:ARG:HG2	1.53	0.72
1:B:581:VAL:HG13	1:B:586:VAL:HB	1.72	0.72
1:D:1520:LEU:CD2	1:D:1611:LEU:HD12	2.16	0.71
1:C:1040:PHE:O	1:C:1044:LYS:HG3	1.92	0.70
1:A:14:HIS:CE1	1:A:16:GLY:HA3	2.27	0.69
1:C:1058:MET:HE1	1:D:1513:LEU:HG	1.74	0.69
1:C:1093:HIS:HB3	1:C:1113:ARG:HH21	1.56	0.69
1:A:57:MET:CE	1:B:515:TYR:HB2	2.25	0.67
1:A:112:LYS:H	1:A:112:LYS:CD	1.92	0.66
1:A:25:ARG:HH12	1:A:29:ASP:C	2.00	0.65
1:C:1019:GLU:O	1:C:1023:ARG:HG3	1.95	0.65
1:C:1077:LYS:O	1:C:1081:VAL:HG23	1.97	0.65
1:D:1525:ARG:NH2	1:D:1529:ASP:HA	2.08	0.65
1:C:1023:ARG:HH22	1:C:1112:LYS:CG	2.07	0.64
1:C:1020:LEU:HD22	1:C:1111:LEU:HD12	1.80	0.64
1:A:37:THR:O	1:A:41:ASN:HB2	1.98	0.64
1:A:62:ILE:HD13	1:B:562:ILE:HD13	1.79	0.63
1:A:14:HIS:O	1:A:16:GLY:N	2.31	0.63
1:C:1017:HIS:CD2	1:C:1017:HIS:N	2.64	0.63
1:D:1520:LEU:HD22	1:D:1611:LEU:CD1	2.20	0.63
1:A:25:ARG:NH1	1:A:25:ARG:HG2	2.12	0.63
1:A:63:ARG:HG2	1:A:63:ARG:HH11	1.64	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1039:GLU:O	1:C:1040:PHE:C	2.35	0.62
1:A:25:ARG:HH11	1:A:25:ARG:HG2	1.63	0.62
1:D:1544:LYS:HG2	1:D:1544:LYS:O	2.00	0.61
1:A:51:ASP:OD2	1:A:54:GLN:HG3	2.00	0.61
1:A:97:GLY:HA2	1:A:100:ASP:OD1	2.00	0.61
1:D:1539:GLU:H	1:D:1539:GLU:CD	2.04	0.61
1:A:23:ARG:HD2	1:A:112:LYS:NZ	2.17	0.60
1:D:1543:ILE:C	1:D:1545:HIS:H	2.04	0.60
1:B:540:PHE:CE2	1:B:544:LYS:HD3	2.37	0.60
1:A:100:ASP:O	1:A:103:LYS:HG3	2.01	0.60
1:B:523:ARG:HG3	1:B:523:ARG:NH1	2.12	0.59
1:C:1051:ASP:OD2	1:C:1054:GLN:HG3	2.02	0.59
1:A:63:ARG:NH1	1:A:63:ARG:HG2	2.17	0.59
1:C:1023:ARG:HH22	1:C:1113:ARG:H	1.49	0.59
1:C:1099:PHE:HA	1:C:1101:PHE:CE1	2.37	0.59
1:C:1094:ASP:H	1:C:1113:ARG:HH22	1.48	0.59
1:A:108:VAL:O	1:A:109:ILE:HD13	2.03	0.59
1:A:41:ASN:HA	1:A:44:LYS:HZ2	1.67	0.58
1:C:1083:LYS:HE3	1:C:1084:PHE:CZ	2.39	0.58
1:C:1012:LEU:HD21	1:C:1049:TYR:HB3	1.86	0.57
1:C:1039:GLU:HA	1:C:1042:GLN:HB3	1.85	0.57
1:C:1057:MET:HE1	1:D:1515:TYR:O	2.04	0.57
1:B:589:PHE:CE1	1:B:591:MET:HG2	2.40	0.57
1:C:1039:GLU:O	1:C:1042:GLN:N	2.38	0.57
1:C:1015:TYR:HB2	1:D:1557:MET:SD	2.45	0.56
1:C:1046:LYS:HG2	1:C:1047:LYS:N	2.19	0.56
1:C:1078:GLU:H	1:C:1078:GLU:CD	2.09	0.55
1:B:503:ARG:NH1	1:B:584:PHE:O	2.37	0.55
1:C:1020:LEU:C	1:C:1020:LEU:HD13	2.26	0.55
1:A:15:TYR:CE1	1:B:557:MET:HB3	2.42	0.55
1:A:14:HIS:HE1	1:A:16:GLY:HA3	1.73	0.54
1:A:15:TYR:HB2	1:B:554:GLN:HG3	1.88	0.54
1:C:1017:HIS:CD2	1:C:1017:HIS:H	2.25	0.54
1:A:93:HIS:HB2	1:A:111:LEU:O	2.08	0.54
1:C:1023:ARG:NH2	1:C:1112:LYS:HG3	2.08	0.54
1:B:588:VAL:HG22	1:B:607:GLU:HB3	1.90	0.54
1:D:1550:TYR:HB3	1:D:1554:GLN:HG3	1.89	0.54
1:A:14:HIS:CD2	1:A:17:HIS:HE1	2.25	0.54
1:A:51:ASP:OD2	1:A:51:ASP:C	2.44	0.54
1:C:1014:HIS:ND1	1:C:1016:GLY:HA3	2.23	0.54
1:C:1112:LYS:HG3	1:C:1113:ARG:N	2.20	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1513:LEU:HA	1:D:1517:HIS:ND1	2.23	0.54
1:B:533:VAL:HG11	1:B:559:LEU:HD13	1.89	0.53
1:C:1039:GLU:C	1:C:1041:ASN:N	2.57	0.53
1:C:1015:TYR:HB2	1:D:1557:MET:CE	2.39	0.53
1:C:1093:HIS:HB3	1:C:1113:ARG:NH2	2.22	0.53
1:C:1014:HIS:HE1	1:C:1016:GLY:HA3	1.69	0.52
1:C:1094:ASP:H	1:C:1113:ARG:NH2	2.06	0.52
1:C:1040:PHE:CD2	1:C:1071:GLU:HG2	2.45	0.52
1:D:1539:GLU:CD	1:D:1539:GLU:N	2.63	0.52
1:A:3:ARG:NH1	1:A:3:ARG:HG2	2.24	0.51
1:C:1056:LYS:O	1:C:1060:GLU:HG3	2.10	0.51
1:C:1007:TYR:CZ	1:C:1077:LYS:HG2	2.46	0.51
1:A:12:LEU:HD21	1:D:1549:TYR:CD1	2.46	0.51
1:A:112:LYS:HG2	1:A:113:ARG:N	2.16	0.51
1:B:520:LEU:HD23	1:B:611:LEU:HD12	1.93	0.51
1:D:1567:LEU:HD12	1:D:1568:VAL:N	2.26	0.51
1:D:1522:ARG:HH22	1:D:1523:ARG:NE	2.10	0.50
1:D:1597:GLY:HA2	1:D:1600:ASP:OD2	2.11	0.50
1:C:1007:TYR:CD2	1:C:1034:ALA:HB3	2.47	0.50
1:C:1041:ASN:O	1:C:1044:LYS:HB2	2.11	0.50
1:C:1044:LYS:O	1:C:1045:HIS:HB2	2.11	0.50
1:A:25:ARG:HH11	1:A:25:ARG:CG	2.25	0.50
1:D:1600:ASP:O	1:D:1603:LYS:HG3	2.12	0.49
1:B:577:LYS:HZ3	1:B:599:PHE:HE1	1.58	0.49
1:C:1112:LYS:O	1:C:1113:ARG:HB2	2.12	0.49
1:A:63:ARG:HG3	1:B:561:SER:O	2.13	0.49
1:D:1593:HIS:HB2	1:D:1611:LEU:O	2.12	0.49
1:B:547:LYS:HE2	1:C:1049:TYR:CD1	2.48	0.48
1:A:23:ARG:HD2	1:A:112:LYS:HZ3	1.78	0.48
1:B:504:VAL:HG12	1:B:505:ILE:N	2.29	0.48
1:B:547:LYS:HE2	1:C:1049:TYR:CE1	2.48	0.48
1:A:19:GLU:OE2	1:A:22:ARG:NH2	2.38	0.48
1:A:35:LEU:HD11	1:A:55:ARG:NH1	2.29	0.48
1:A:57:MET:HE2	1:B:515:TYR:HB2	1.96	0.48
1:A:2:LYS:HA	1:A:87:ASP:OD2	2.14	0.48
1:C:1002:LYS:N	1:C:1029:ASP:OD2	2.41	0.47
1:C:1039:GLU:O	1:C:1041:ASN:N	2.47	0.47
1:D:1535:LEU:HD12	1:D:1552:TYR:CD1	2.49	0.47
1:C:1054:GLN:HB3	1:D:1515:TYR:CD2	2.50	0.47
1:A:44:LYS:HB2	1:A:44:LYS:NZ	2.30	0.47
1:B:520:LEU:HD13	1:B:521:LEU:N	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:ARG:CG	1:A:3:ARG:HH11	2.24	0.46
1:A:37:THR:HG21	1:A:72:LYS:HG2	1.96	0.46
1:B:523:ARG:CG	1:B:523:ARG:NH1	2.75	0.46
1:C:1023:ARG:NH2	1:C:1113:ARG:H	2.13	0.46
1:A:40:PHE:CE2	1:A:44:LYS:HE3	2.50	0.46
1:C:1014:HIS:ND1	1:C:1016:GLY:N	2.64	0.46
1:A:112:LYS:O	1:C:1094:ASP:OD1	2.33	0.46
1:D:1546:LYS:HG3	1:D:1546:LYS:O	2.15	0.46
1:B:507:TYR:CD2	1:B:534:ALA:HB3	2.51	0.46
1:B:577:LYS:O	1:B:581:VAL:HG23	2.16	0.46
1:C:1023:ARG:HE	1:C:1111:LEU:HD13	1.81	0.46
1:D:1501:MET:HG2	1:D:1503:ARG:CZ	2.46	0.45
1:B:530:TYR:HA	1:B:566:ASP:OD2	2.16	0.45
1:C:1097:GLY:HA2	1:C:1100:ASP:OD1	2.16	0.45
1:A:15:TYR:H	1:B:554:GLN:HE21	1.64	0.45
1:D:1513:LEU:HA	1:D:1517:HIS:CE1	2.52	0.45
1:C:1014:HIS:C	1:C:1016:GLY:N	2.69	0.45
1:B:540:PHE:O	1:B:544:LYS:HG3	2.16	0.45
1:B:593:HIS:HB2	1:B:611:LEU:O	2.16	0.45
1:A:10:TYR:HB3	1:A:13:LEU:HD21	1.99	0.45
1:C:1012:LEU:HD21	1:C:1049:TYR:CB	2.46	0.45
1:B:563:ARG:HD2	1:B:564:TYR:CZ	2.53	0.44
1:B:587:ASP:OD2	1:B:587:ASP:N	2.51	0.44
1:A:95:TRP:CE3	1:A:99:PHE:CE2	3.05	0.44
1:C:1051:ASP:OD2	1:C:1051:ASP:C	2.56	0.44
1:A:35:LEU:HD11	1:A:55:ARG:HH11	1.83	0.44
1:B:522:ARG:HG3	1:B:564:TYR:CE1	2.53	0.44
1:C:1052:TYR:CA	1:C:1055:ARG:HH21	2.31	0.43
1:D:1543:ILE:C	1:D:1545:HIS:N	2.70	0.43
1:B:593:HIS:HB2	1:B:612:LYS:O	2.18	0.43
1:A:63:ARG:NH1	1:B:563:ARG:HA	2.33	0.43
1:B:552:TYR:OH	1:B:570:PRO:HB3	2.18	0.43
1:A:1:MET:HE2	1:A:3:ARG:HH22	1.84	0.43
1:B:613:ARG:HG2	1:B:613:ARG:HH11	1.83	0.43
1:C:1039:GLU:HA	1:C:1042:GLN:CB	2.46	0.43
1:C:1003:ARG:NH1	1:C:1084:PHE:O	2.48	0.43
1:B:509:THR:HG23	1:B:555:ARG:NH1	2.33	0.43
1:C:1017:HIS:HD2	1:C:1017:HIS:N	2.11	0.43
1:D:1525:ARG:HH21	1:D:1529:ASP:CA	2.14	0.43
1:D:1520:LEU:C	1:D:1520:LEU:HD13	2.39	0.43
1:C:1005:ILE:HG22	1:C:1086:VAL:HG11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:HIS:C	1:A:16:GLY:N	2.72	0.43
1:B:522:ARG:NH1	1:B:526:GLU:OE2	2.52	0.42
1:C:1078:GLU:N	1:C:1078:GLU:OE2	2.52	0.42
1:D:1596:GLU:HA	1:D:1610:TYR:CD1	2.54	0.42
1:B:540:PHE:CZ	1:B:544:LYS:HD3	2.53	0.42
1:C:1038:ASP:O	1:C:1042:GLN:N	2.43	0.42
1:D:1554:GLN:O	1:D:1557:MET:HG3	2.19	0.42
1:C:1005:ILE:CG2	1:C:1086:VAL:HG11	2.50	0.42
1:D:1522:ARG:HH22	1:D:1523:ARG:HE	1.65	0.42
1:C:1046:LYS:CG	1:C:1047:LYS:N	2.83	0.42
1:D:1581:VAL:HG13	1:D:1586:VAL:HB	2.01	0.42
1:A:25:ARG:HG3	1:A:31:LEU:CB	2.50	0.42
1:B:520:LEU:HD13	1:B:520:LEU:C	2.40	0.42
1:C:1022:ARG:HG2	1:C:1026:GLU:OE2	2.20	0.42
1:A:14:HIS:CD2	1:A:17:HIS:CE1	3.06	0.42
1:D:1557:MET:HG3	1:D:1558:MET:N	2.34	0.42
1:A:57:MET:HE1	1:B:515:TYR:HB2	2.01	0.41
1:C:1040:PHE:CE1	1:C:1073:GLY:HA2	2.55	0.41
1:A:39:GLU:N	1:A:39:GLU:CD	2.73	0.41
1:A:35:LEU:HD13	1:A:59:LEU:HD11	2.01	0.41
1:D:1599:PHE:HA	1:D:1601:PHE:CE1	2.55	0.41
1:B:509:THR:O	1:B:535:LEU:HD11	2.20	0.41
1:C:1052:TYR:HA	1:C:1055:ARG:NH2	2.36	0.41
1:B:535:LEU:HD21	1:B:555:ARG:HB3	2.03	0.41
1:A:81:VAL:HA	1:A:86:VAL:HB	2.03	0.41
1:B:596:GLU:HA	1:B:610:TYR:CD1	2.55	0.41
1:D:1567:LEU:HD12	1:D:1568:VAL:H	1.85	0.41
1:A:3:ARG:CG	1:A:3:ARG:NH1	2.80	0.40
1:B:554:GLN:O	1:B:558:MET:HG3	2.20	0.40
1:C:1014:HIS:O	1:C:1016:GLY:N	2.54	0.40
1:A:62:ILE:HD13	1:B:562:ILE:CD1	2.48	0.40
1:D:1501:MET:CE	1:D:1503:ARG:NH2	2.83	0.40
1:D:1551:ASP:H	1:D:1554:GLN:HE21	1.69	0.40
1:A:100:ASP:O	1:A:103:LYS:CG	2.68	0.40
1:D:1577:LYS:O	1:D:1581:VAL:HG23	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	113/132 (86%)	98 (87%)	13 (12%)	2 (2%)	10	43
1	B	111/132 (84%)	104 (94%)	6 (5%)	1 (1%)	20	62
1	C	111/132 (84%)	102 (92%)	9 (8%)	0	100	100
1	D	111/132 (84%)	100 (90%)	9 (8%)	2 (2%)	10	43
All	All	446/528 (84%)	404 (91%)	37 (8%)	5 (1%)	17	56

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	15	TYR
1	A	113	ARG
1	B	612	LYS
1	D	1612	LYS
1	D	1544	LYS

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	106/120 (88%)	101 (95%)	5 (5%)	30	69
1	B	104/120 (87%)	96 (92%)	8 (8%)	15	48
1	C	104/120 (87%)	101 (97%)	3 (3%)	48	82
1	D	104/120 (87%)	94 (90%)	10 (10%)	10	36
All	All	418/480 (87%)	392 (94%)	26 (6%)	21	58

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ARG
1	A	41	ASN
1	A	79	ASP
1	A	100	ASP
1	A	112	LYS
1	B	513	LEU
1	B	520	LEU
1	B	523	ARG
1	B	541	ASN
1	B	551	ASP
1	B	579	ASP
1	B	587	ASP
1	B	604	ASP
1	C	1003	ARG
1	C	1035	LEU
1	C	1100	ASP
1	D	1501	MET
1	D	1503	ARG
1	D	1519	GLU
1	D	1525	ARG
1	D	1535	LEU
1	D	1541	ASN
1	D	1557	MET
1	D	1582	GLU
1	D	1583	LYS
1	D	1604	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	B	554	GLN
1	C	1041	ASN
1	C	1093	HIS
1	D	1554	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 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.