



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:01 PM UTC

PDB ID : 4IZI / pdb_00004izi
Title : Crystal Structure of the Alpha1 dimer of Thermus thermophilus Transhydrogenase in P4(3)
Authors : Stout, C.D.; Yamaguchi, M.; Leung, J.
Deposited on : 2013-01-30
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

<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
Mogul : 2022.3.0, CSD as543be (2022)
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

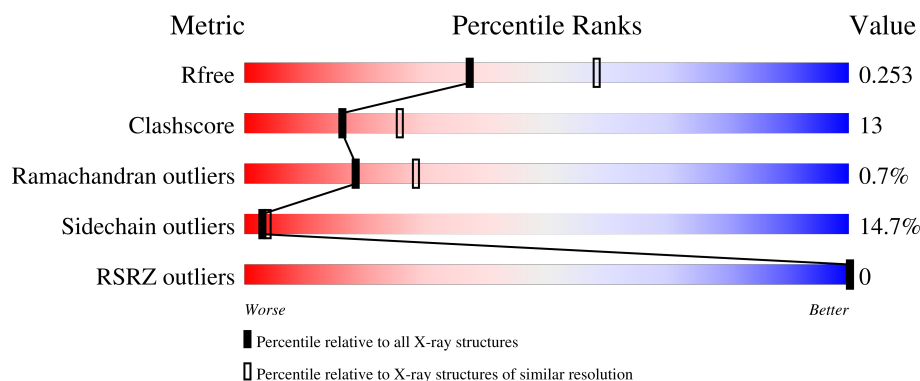
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (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 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	381	 69% 20% 6% .
1	B	381	 71% 18% 6% . .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	A	403	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5729 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 NAD/NADP transhydrogenase alpha subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	364	Total	C	N	O	S	0	1	0
			2745	1743	490	501	11			
1	B	366	Total	C	N	O	S	0	1	0
			2755	1749	492	503	11			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	expression tag	UNP Q72GR8
A	-4	HIS	-	expression tag	UNP Q72GR8
A	-3	HIS	-	expression tag	UNP Q72GR8
A	-2	HIS	-	expression tag	UNP Q72GR8
A	-1	HIS	-	expression tag	UNP Q72GR8
A	0	HIS	-	expression tag	UNP Q72GR8
B	-5	HIS	-	expression tag	UNP Q72GR8
B	-4	HIS	-	expression tag	UNP Q72GR8
B	-3	HIS	-	expression tag	UNP Q72GR8
B	-2	HIS	-	expression tag	UNP Q72GR8
B	-1	HIS	-	expression tag	UNP Q72GR8
B	0	HIS	-	expression tag	UNP Q72GR8

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		

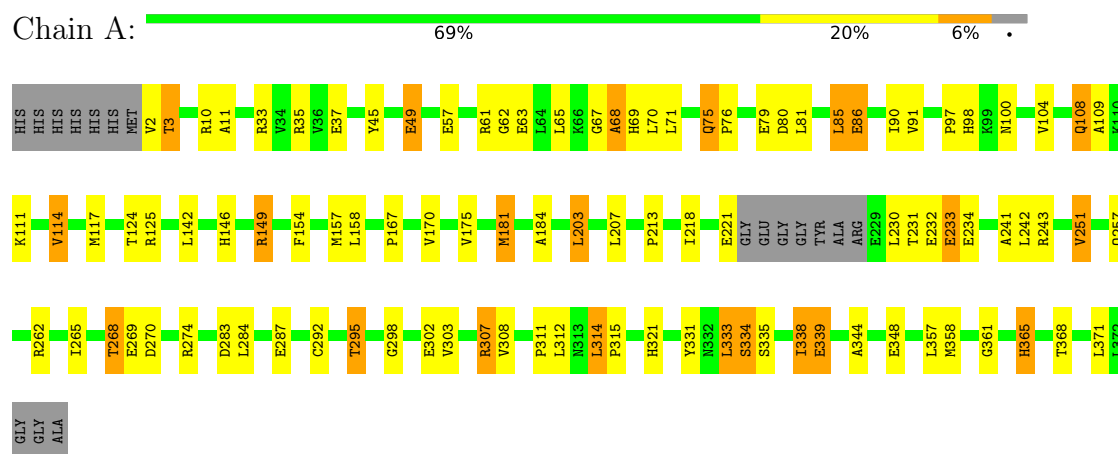
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	90	Total	O	0	0
			90	90		
3	B	97	Total	O	0	0
			97	97		

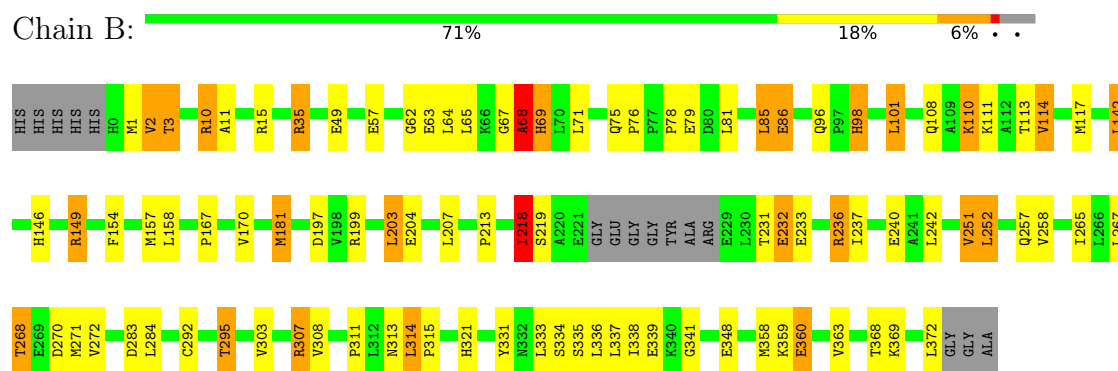
3 Residue-property plots

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: NAD/NADP transhydrogenase alpha subunit 1



- Molecule 1: NAD/NADP transhydrogenase alpha subunit 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	85.79Å 85.79Å 146.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.37 – 2.40 38.37 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (38.37-2.40) 99.8 (38.37-2.40)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.202 , 0.257 0.197 , 0.253	Depositor DCC
R_{free} test set	2087 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	59.0	Xtriage
Anisotropy	0.008	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.487 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5729	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.93% 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: GOL

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	0.95	1/2792 (0.0%)	1.03	3/3787 (0.1%)
1	B	0.97	2/2802 (0.1%)	1.06	6/3801 (0.2%)
All	All	0.96	3/5594 (0.1%)	1.05	9/7588 (0.1%)

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	2
1	B	0	1
All	All	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	258	VAL	CA-CB	5.84	1.58	1.54
1	B	1	MET	CA-C	5.43	1.58	1.52
1	A	241	ALA	CA-C	5.06	1.59	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	68	ALA	N-CA-C	9.95	125.25	111.92
1	B	2	VAL	N-CA-C	7.38	120.26	109.63
1	A	63	GLU	N-CA-C	-7.15	104.37	113.23
1	A	85	LEU	N-CA-C	6.98	119.53	111.02
1	B	68	ALA	N-CA-C	6.25	124.11	110.80
1	B	218	ILE	N-CA-CB	5.70	116.93	110.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	69	HIS	N-CA-C	5.46	120.07	113.41
1	B	63	GLU	N-CA-C	-5.43	106.66	113.28
1	B	85	LEU	N-CA-C	5.28	117.46	111.02

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	62	GLY	Peptide
1	A	67	GLY	Peptide
1	B	68	ALA	Peptide

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	2745	0	2869	76	0
1	B	2755	0	2876	80	0
2	A	30	0	40	6	0
2	B	12	0	16	3	0
3	A	90	0	0	4	0
3	B	97	0	0	12	0
All	All	5729	0	5801	149	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 13.

All (149) 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:146:HIS:ND1	1:B:149[A]:ARG:NH1	1.88	1.21
1:A:312:LEU:H	2:A:403:GOL:H11	1.20	1.04
1:A:146:HIS:ND1	1:A:149[A]:ARG:NH1	2.07	1.03
1:B:181:MET:HE2	1:B:181:MET:HA	1.40	1.03
1:A:65:LEU:O	1:A:68:ALA:HB2	1.67	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:PHE:H	1:B:321:HIS:HD2	1.17	0.92
1:A:312:LEU:H	2:A:403:GOL:C1	1.85	0.90
1:B:203:LEU:HD21	2:B:402:GOL:H11	1.53	0.89
1:B:181:MET:HE3	2:B:401:GOL:H2	1.53	0.88
1:A:321:HIS:HD2	1:B:154:PHE:H	1.24	0.86
1:B:313:ASN:HB2	3:B:565:HOH:O	1.75	0.86
1:A:149[A]:ARG:HD2	1:B:149[A]:ARG:HD2	1.61	0.83
1:A:231:THR:HG23	1:A:234:GLU:H	1.41	0.83
1:B:110:LYS:HE3	3:B:579:HOH:O	1.78	0.82
1:B:98:HIS:H	1:B:98:HIS:CD2	1.99	0.81
1:B:268:THR:CG2	1:B:270:ASP:OD1	2.28	0.81
1:B:62:GLY:HA2	3:B:561:HOH:O	1.81	0.80
1:A:292:CYS:HB3	1:A:295:THR:CG2	2.11	0.80
1:A:231:THR:CG2	1:A:234:GLU:HG3	2.12	0.80
1:B:313:ASN:CB	3:B:565:HOH:O	2.29	0.79
1:B:181:MET:HA	1:B:181:MET:CE	2.14	0.78
1:A:268:THR:CG2	1:A:270:ASP:OD1	2.31	0.77
1:B:114:VAL:CG1	1:B:358:MET:HG2	2.15	0.77
1:B:292:CYS:HB3	1:B:295:THR:CG2	2.14	0.77
1:B:65:LEU:HD23	1:B:85:LEU:HA	1.68	0.75
1:A:312:LEU:N	2:A:403:GOL:H11	2.00	0.75
1:A:311:PRO:HB2	1:A:314:LEU:HD13	1.70	0.74
1:B:331:TYR:O	1:B:334:SER:HB2	1.88	0.74
1:A:181:MET:HA	1:A:181:MET:HE2	1.69	0.74
1:A:269:GLU:HG3	3:A:554:HOH:O	1.87	0.74
1:A:274:ARG:HD3	3:A:576:HOH:O	1.88	0.73
1:B:114:VAL:HG12	1:B:358:MET:HG2	1.70	0.73
1:B:67:GLY:O	1:B:68:ALA:HB2	1.90	0.72
1:B:85:LEU:O	1:B:86:GLU:HB3	1.89	0.72
1:A:243:ARG:HD2	1:A:274:ARG:HH11	1.55	0.71
1:A:114:VAL:HG13	1:A:358:MET:HG2	1.72	0.70
1:A:181:MET:HE3	2:A:401:GOL:H2	1.74	0.70
1:B:283:ASP:OD2	1:B:295:THR:HG21	1.92	0.69
1:A:283:ASP:OD2	1:A:295:THR:HG21	1.92	0.69
1:B:3:THR:HG23	1:B:68:ALA:HB3	1.75	0.69
1:A:181:MET:HE2	1:A:184:ALA:HB3	1.74	0.69
1:B:170:VAL:HG13	1:B:251:VAL:HG22	1.75	0.67
1:A:243:ARG:HD2	1:A:274:ARG:NH1	2.11	0.64
1:B:311:PRO:HB2	1:B:314:LEU:HD13	1.79	0.64
1:A:233:GLU:OE1	1:A:233:GLU:HA	1.98	0.63
1:B:117:MET:HA	1:B:117:MET:HE2	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:292:CYS:HB3	1:B:295:THR:HG23	1.81	0.63
1:A:100:ASN:O	1:A:104:VAL:HG23	1.99	0.62
1:A:65:LEU:O	1:A:68:ALA:CB	2.47	0.62
1:A:181:MET:HA	1:A:181:MET:CE	2.29	0.62
1:B:363:VAL:O	1:B:369:LYS:HE3	1.99	0.62
1:A:298:GLY:HA2	2:A:403:GOL:H12	1.82	0.62
1:A:231:THR:HG22	1:A:234:GLU:HG3	1.83	0.61
1:B:268:THR:HG21	1:B:270:ASP:OD1	2.00	0.61
1:B:149[A]:ARG:NH2	3:B:587:HOH:O	2.31	0.61
1:B:114:VAL:HG12	1:B:358:MET:CG	2.31	0.61
1:A:157:MET:HB2	1:A:167:PRO:HD3	1.83	0.60
1:A:146:HIS:CE1	1:A:149[A]:ARG:NH1	2.69	0.60
1:A:154:PHE:H	1:B:321:HIS:CD2	2.08	0.60
1:A:203:LEU:HG	1:A:213:PRO:HB3	1.83	0.60
1:A:85:LEU:O	1:A:86:GLU:CB	2.51	0.59
1:B:236:ARG:O	1:B:240:GLU:HG3	2.03	0.59
1:A:331:TYR:O	1:A:334:SER:HB2	2.03	0.59
1:A:79:GLU:OE1	3:A:550:HOH:O	2.16	0.59
1:A:98:HIS:H	1:A:98:HIS:CD2	2.19	0.59
1:B:218:ILE:HD12	1:B:219:SER:N	2.19	0.58
1:B:85:LEU:O	1:B:86:GLU:CB	2.50	0.58
1:B:98:HIS:CE1	3:B:586:HOH:O	2.55	0.58
1:A:268:THR:HG21	1:A:270:ASP:OD1	2.03	0.58
1:A:70:LEU:CD1	1:A:338:ILE:HD11	2.34	0.57
1:B:157:MET:HB2	1:B:167:PRO:HD3	1.86	0.57
1:A:70:LEU:HD12	1:A:338:ILE:HD11	1.86	0.57
1:A:231:THR:HG22	1:A:234:GLU:CG	2.35	0.57
1:B:295:THR:HG22	1:B:308:VAL:HG11	1.87	0.56
1:B:113:THR:OG1	1:B:359:LYS:HA	2.05	0.56
1:B:68:ALA:HB1	1:B:69:HIS:CG	2.41	0.56
1:A:97:PRO:HA	1:A:104:VAL:HG21	1.88	0.55
1:B:146:HIS:CE1	1:B:149[A]:ARG:NH1	2.69	0.55
1:B:35:ARG:HD3	1:B:64:LEU:O	2.06	0.54
1:A:365:HIS:CD2	1:A:368:THR:OG1	2.61	0.54
1:B:295:THR:HG22	1:B:308:VAL:CG1	2.37	0.54
1:B:197:ASP:O	2:B:402:GOL:H12	2.06	0.54
1:A:321:HIS:HD2	1:B:154:PHE:N	1.99	0.53
1:B:170:VAL:HG13	1:B:251:VAL:CG2	2.39	0.53
1:A:170:VAL:HG13	1:A:251:VAL:HG22	1.91	0.52
1:B:35:ARG:HG3	1:B:64:LEU:HD12	1.92	0.51
1:B:2:VAL:HG11	1:B:338:ILE:HG12	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:334:SER:O	1:B:336:LEU:N	2.43	0.51
1:A:292:CYS:HB3	1:A:295:THR:HG23	1.90	0.51
1:A:314:LEU:N	1:A:315:PRO:CD	2.74	0.51
1:A:70:LEU:HD23	1:A:90:ILE:HB	1.91	0.51
1:A:75:GLN:HB3	1:A:76:PRO:HD2	1.93	0.51
1:A:292:CYS:O	1:A:295:THR:HG23	2.10	0.51
1:B:10:ARG:O	1:B:11:ALA:C	2.53	0.51
1:B:314:LEU:N	1:B:315:PRO:CD	2.75	0.50
1:A:312:LEU:O	2:A:403:GOL:H11	2.12	0.50
1:A:117:MET:HA	1:A:117:MET:HE2	1.93	0.50
1:B:75:GLN:HB3	1:B:76:PRO:HD2	1.94	0.50
1:A:10:ARG:O	1:A:11:ALA:C	2.55	0.49
1:B:203:LEU:HG	1:B:213:PRO:HB3	1.93	0.49
1:A:339:GLU:HB2	1:A:344:ALA:HB3	1.93	0.49
1:A:3:THR:H	1:A:69:HIS:HD2	1.60	0.49
1:B:3:THR:HG23	1:B:68:ALA:CB	2.42	0.49
1:B:67:GLY:O	1:B:68:ALA:CB	2.57	0.49
1:A:37:GLU:OE1	1:A:61:ARG:NH2	2.45	0.49
1:A:2:VAL:HG11	1:A:338:ILE:HG12	1.95	0.49
1:B:232:GLU:HG3	3:B:597:HOH:O	2.12	0.49
1:B:233:GLU:HA	1:B:233:GLU:OE1	2.13	0.48
1:A:80:ASP:OD1	1:A:81:LEU:N	2.46	0.48
1:B:334:SER:C	1:B:336:LEU:H	2.22	0.48
1:B:232:GLU:HB3	3:B:597:HOH:O	2.12	0.48
1:B:368:THR:O	1:B:372:LEU:HB2	2.15	0.47
1:A:91:VAL:HB	1:A:114:VAL:HB	1.97	0.46
1:A:65:LEU:HD23	1:A:85:LEU:HA	1.96	0.46
1:A:154:PHE:N	1:B:321:HIS:HD2	1.99	0.46
1:A:302:GLU:OE2	1:A:307:ARG:HD3	2.15	0.46
1:A:295:THR:HG22	1:A:308:VAL:HG11	1.98	0.46
1:B:334:SER:C	1:B:336:LEU:N	2.72	0.46
1:A:175:VAL:HG12	1:A:175:VAL:O	2.17	0.45
1:A:365:HIS:HD2	1:A:368:THR:OG1	1.97	0.45
1:B:68:ALA:HB1	1:B:69:HIS:CD2	2.51	0.45
1:A:274:ARG:CD	3:A:576:HOH:O	2.59	0.45
1:A:231:THR:CG2	1:A:234:GLU:CG	2.89	0.45
1:B:15:ARG:HD2	3:B:546:HOH:O	2.16	0.45
1:A:45:TYR:OH	1:B:307:ARG:NH1	2.50	0.45
1:B:267:LEU:HD23	1:B:267:LEU:HA	1.76	0.44
1:B:359:LYS:O	1:B:360:GLU:C	2.61	0.43
1:A:114:VAL:HG13	1:A:358:MET:CG	2.45	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:LEU:HG	1:B:65:LEU:HD12	2.00	0.43
1:B:2:VAL:N	1:B:341:GLY:O	2.52	0.43
1:A:70:LEU:CD2	1:A:90:ILE:HB	2.48	0.43
1:A:111:LYS:HA	1:A:361:GLY:HA3	2.00	0.43
1:A:108:GLN:HG3	1:A:109:ALA:N	2.34	0.42
1:B:101:LEU:HD23	1:B:101:LEU:HA	1.88	0.42
1:B:252:LEU:HD22	1:B:252:LEU:N	2.34	0.42
1:B:2:VAL:HG22	1:B:3:THR:N	2.35	0.42
1:B:78:PRO:HD2	1:B:81:LEU:HD12	2.02	0.42
1:A:117:MET:HE1	1:A:333:LEU:HD11	2.02	0.42
1:A:3:THR:HA	1:A:33:ARG:O	2.20	0.41
1:B:79:GLU:HB2	3:B:592:HOH:O	2.19	0.41
1:A:117:MET:HE1	1:A:333:LEU:CD1	2.50	0.41
1:A:49:GLU:H	1:A:49:GLU:CD	2.28	0.41
1:B:232:GLU:CB	3:B:597:HOH:O	2.68	0.41
1:B:337:LEU:C	1:B:338:ILE:HG13	2.46	0.41
1:B:142:LEU:HD12	1:B:142:LEU:HA	1.95	0.41
1:A:231:THR:HG22	1:A:234:GLU:CD	2.46	0.41
1:B:268:THR:HG22	1:B:271:MET:H	1.85	0.40
1:B:292:CYS:O	1:B:295:THR:HG23	2.21	0.40
1:B:218:ILE:HD11	3:B:539:HOH:O	2.21	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	361/381 (95%)	347 (96%)	12 (3%)	2 (1%)	21	32
1	B	363/381 (95%)	344 (95%)	16 (4%)	3 (1%)	16	25
All	All	724/762 (95%)	691 (95%)	28 (4%)	5 (1%)	18	28

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	68	ALA
1	B	335	SER
1	A	86	GLU
1	A	335	SER
1	B	86	GLU

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	284/293 (97%)	242 (85%)	42 (15%)	3	4
1	B	284/293 (97%)	241 (85%)	43 (15%)	3	3
All	All	568/586 (97%)	483 (85%)	85 (15%)	3	3

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	THR
1	A	35	ARG
1	A	49	GLU
1	A	57	GLU
1	A	71	LEU
1	A	75	GLN
1	A	108	GLN
1	A	114	VAL
1	A	124	THR
1	A	125	ARG
1	A	142	LEU
1	A	149[A]	ARG
1	A	149[B]	ARG
1	A	158	LEU
1	A	181	MET
1	A	203	LEU
1	A	207	LEU
1	A	218	ILE

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Mol	Chain	Res	Type
1	A	221	GLU
1	A	230	LEU
1	A	232	GLU
1	A	233	GLU
1	A	242	LEU
1	A	251	VAL
1	A	257	GLN
1	A	262	ARG
1	A	265	ILE
1	A	268	THR
1	A	284	LEU
1	A	287	GLU
1	A	295	THR
1	A	303	VAL
1	A	307	ARG
1	A	314	LEU
1	A	333	LEU
1	A	334	SER
1	A	338	ILE
1	A	339	GLU
1	A	348	GLU
1	A	357	LEU
1	A	365	HIS
1	A	371	LEU
1	B	3	THR
1	B	10	ARG
1	B	35	ARG
1	B	49	GLU
1	B	57	GLU
1	B	71	LEU
1	B	96	GLN
1	B	98	HIS
1	B	101	LEU
1	B	108	GLN
1	B	110	LYS
1	B	111	LYS
1	B	114	VAL
1	B	142	LEU
1	B	149[A]	ARG
1	B	149[B]	ARG
1	B	158	LEU
1	B	181	MET

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Mol	Chain	Res	Type
1	B	199	ARG
1	B	203	LEU
1	B	204	GLU
1	B	207	LEU
1	B	218	ILE
1	B	231	THR
1	B	232	GLU
1	B	236	ARG
1	B	237	ILE
1	B	242	LEU
1	B	251	VAL
1	B	252	LEU
1	B	257	GLN
1	B	265	ILE
1	B	268	THR
1	B	272	VAL
1	B	284	LEU
1	B	295	THR
1	B	303	VAL
1	B	307	ARG
1	B	314	LEU
1	B	333	LEU
1	B	339	GLU
1	B	348	GLU
1	B	360	GLU

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	52	GLN
1	A	69	HIS
1	A	98	HIS
1	A	108	GLN
1	A	205	GLN
1	A	321	HIS
1	A	332	ASN
1	A	365	HIS
1	B	69	HIS
1	B	98	HIS
1	B	108	GLN
1	B	205	GLN
1	B	321	HIS

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Mol	Chain	Res	Type
1	B	332	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GOL	A	405	-	5,5,5	0.50	0	5,5,5	1.03	0
2	GOL	A	402	-	5,5,5	0.39	0	5,5,5	0.58	0
2	GOL	A	404	-	5,5,5	0.72	0	5,5,5	1.01	0
2	GOL	A	403	-	5,5,5	0.80	0	5,5,5	1.04	0
2	GOL	A	401	-	5,5,5	0.41	0	5,5,5	0.61	0
2	GOL	B	401	-	5,5,5	0.41	0	5,5,5	0.62	0
2	GOL	B	402	-	5,5,5	0.41	0	5,5,5	0.58	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	405	-	-	2/4/4/4	-
2	GOL	A	402	-	-	2/4/4/4	-
2	GOL	A	404	-	-	1/4/4/4	-
2	GOL	A	403	-	-	1/4/4/4	-
2	GOL	A	401	-	-	2/4/4/4	-
2	GOL	B	401	-	-	2/4/4/4	-
2	GOL	B	402	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	GOL	C1-C2-C3-O3
2	A	402	GOL	O1-C1-C2-C3
2	A	405	GOL	O1-C1-C2-C3
2	B	401	GOL	C1-C2-C3-O3
2	B	401	GOL	O2-C2-C3-O3
2	B	402	GOL	O1-C1-C2-C3
2	A	401	GOL	O2-C2-C3-O3
2	A	402	GOL	O1-C1-C2-O2
2	B	402	GOL	O1-C1-C2-O2
2	A	405	GOL	O1-C1-C2-O2
2	A	403	GOL	C1-C2-C3-O3
2	B	402	GOL	O2-C2-C3-O3
2	A	404	GOL	O2-C2-C3-O3

There are no ring outliers.

4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	403	GOL	5	0
2	A	401	GOL	1	0
2	B	401	GOL	1	0
2	B	402	GOL	2	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	364/381 (95%)	-1.43	0 100 100	22, 60, 92, 98	1 (0%)
1	B	366/381 (96%)	-1.44	0 100 100	21, 60, 92, 98	1 (0%)
All	All	730/762 (95%)	-1.44	0 100 100	21, 60, 92, 98	2 (0%)

There are no RSRZ outliers to report.

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(Å ²)	Q<0.9
2	GOL	A	403	6/6	0.98	0.06	41,45,46,53	0
2	GOL	A	405	6/6	0.98	0.07	48,50,56,56	0
2	GOL	A	401	6/6	0.99	0.06	64,67,68,70	0
2	GOL	A	402	6/6	0.99	0.04	82,86,86,87	0
2	GOL	B	401	6/6	0.99	0.06	66,68,70,70	0
2	GOL	B	402	6/6	0.99	0.05	93,96,97,99	0
2	GOL	A	404	6/6	1.00	0.03	48,51,52,60	0

6.5 Other polymers ⓘ

There are no such residues in this entry.