



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:20 PM UTC

PDB ID : 4UNK / pdb_00004unk
Title : Crystal structure of human triosephosphate isomerase (mutant N15D)
Authors : DeLaMora-DeLaMora, I.; Torres-Larios, A.; Enriquez-Flores, S.; Mendez, S.T.; Castillo-Villanueva, A.; Gomez-Manzo, S.; Lopez-Velazquez, G.; Marcial-Quino, J.; Torres-Arroyo, A.; Garcia-Torres, I.; Reyes-Vivas, H.; Oria-Hernandez, J.
Deposited on : 2014-05-29
Resolution : 2.00 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

<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

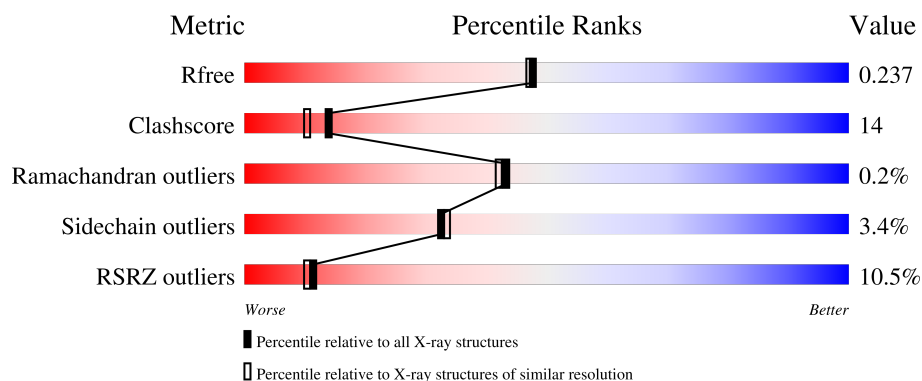
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
R_{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 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	250	7% 64% 31% ••
1	B	250	13% 66% 25% 5% •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4122 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 TRIOSEPHOSPHATE ISOMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total	C	N	O	S	0	0	0
			1844	1167	318	352	7			
1	B	242	Total	C	N	O	S	0	0	0
			1826	1156	315	348	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P60174
A	0	SER	-	expression tag	UNP P60174
A	15	ASP	ASN	engineered mutation	UNP P60174
B	-1	GLY	-	expression tag	UNP P60174
B	0	SER	-	expression tag	UNP P60174
B	15	ASP	ASN	engineered mutation	UNP P60174

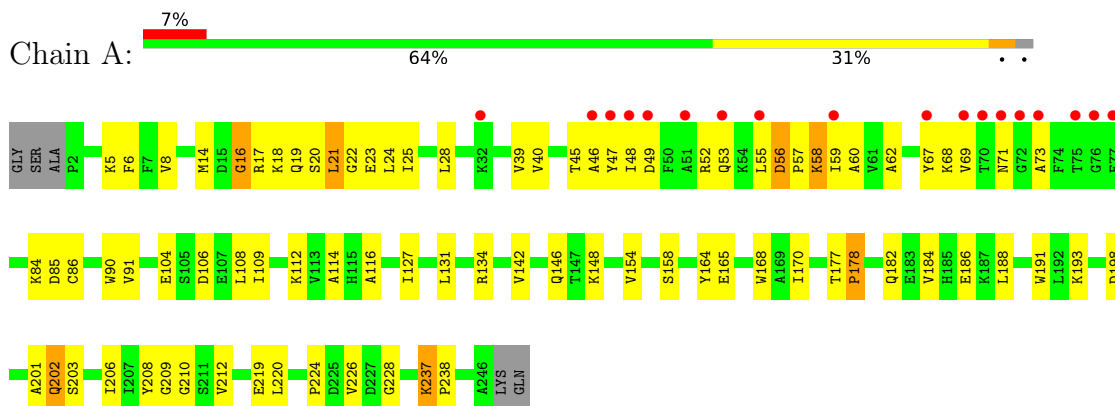
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	304	Total	O	0	0
			304	304		
2	B	148	Total	O	0	0
			148	148		

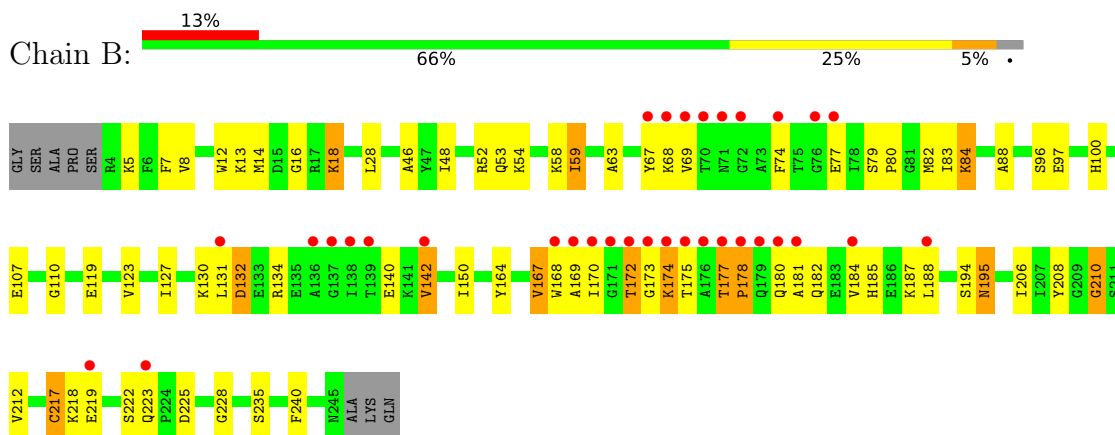
3 Residue-property plots

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: TRIOSEPHOSPHATE ISOMERASE



• Molecule 1: TRIOSEPHOSPHATE ISOMERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	47.63Å 70.71Å 146.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.29 – 2.00 45.29 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.4 (45.29-2.00) 99.4 (45.29-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.08 (at 2.00Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.179 , 0.224 (Not available) , 0.237	Depositor DCC
R_{free} test set	1723 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	22.9	Xtriage
Anisotropy	0.523	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 53.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4122	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 5.07% 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	1.78	42/1879 (2.2%)	1.08	6/2546 (0.2%)
1	B	1.61	14/1860 (0.8%)	1.11	10/2520 (0.4%)
All	All	1.70	56/3739 (1.5%)	1.10	16/5066 (0.3%)

All (56) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	108	LEU	C-O	-7.03	1.15	1.24
1	B	52	ARG	C-O	-6.82	1.16	1.24
1	A	148	LYS	C-O	-6.63	1.16	1.24
1	A	158	SER	C-O	-6.62	1.16	1.24
1	A	20	SER	C-O	-6.29	1.16	1.24
1	A	28	LEU	C-O	-6.27	1.16	1.24
1	A	209	GLY	C-O	-6.22	1.18	1.24
1	A	237	LYS	C-O	-6.11	1.17	1.24
1	A	23	GLU	C-O	-6.00	1.17	1.24
1	A	84	LYS	C-O	-5.84	1.17	1.24
1	A	24	LEU	C-O	-5.82	1.17	1.24
1	A	131	LEU	C-O	-5.81	1.17	1.24
1	A	142	VAL	C-O	-5.80	1.17	1.24
1	B	88	ALA	C-O	-5.76	1.16	1.23
1	B	46	ALA	C-O	-5.73	1.17	1.24
1	A	201	ALA	C-O	-5.72	1.17	1.24
1	A	202	GLN	C-O	-5.66	1.17	1.24
1	B	18	LYS	C-O	-5.65	1.17	1.24
1	A	114	ALA	C-O	-5.64	1.17	1.24
1	A	226	VAL	C-O	-5.63	1.17	1.24
1	A	22	GLY	C-O	-5.59	1.17	1.23
1	B	150	ILE	C-O	-5.58	1.17	1.24
1	A	19	GLN	C-O	-5.56	1.17	1.24
1	A	8	VAL	C-O	-5.52	1.18	1.24
1	A	191	TRP	C-O	-5.52	1.17	1.24
1	A	18	LYS	C-O	-5.52	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	165	GLU	C-O	-5.52	1.18	1.24
1	A	203	SER	C-O	-5.51	1.16	1.24
1	A	188	LEU	C-O	-5.48	1.17	1.24
1	A	39	VAL	C-O	-5.46	1.18	1.24
1	B	142	VAL	C-O	-5.45	1.18	1.24
1	B	110	GLY	C-O	-5.45	1.17	1.23
1	B	53	GLN	C-O	-5.42	1.17	1.24
1	A	116	ALA	C-O	-5.40	1.17	1.24
1	A	109	ILE	C-O	-5.38	1.17	1.24
1	A	212	VAL	C-O	-5.38	1.18	1.24
1	A	57	PRO	N-CD	5.34	1.55	1.47
1	A	178	PRO	N-CD	5.32	1.55	1.47
1	B	178	PRO	N-CD	5.30	1.55	1.47
1	B	8	VAL	C-O	-5.26	1.18	1.24
1	A	91	VAL	C-O	-5.23	1.18	1.24
1	B	48	ILE	C-O	-5.22	1.18	1.24
1	A	85	ASP	C-O	-5.19	1.17	1.24
1	A	146	GLN	C-O	-5.17	1.18	1.24
1	A	86	CYS	C-O	-5.13	1.17	1.24
1	B	177	THR	C-N	5.13	1.41	1.33
1	A	5	LYS	C-O	-5.12	1.17	1.23
1	B	123	VAL	C-O	-5.12	1.18	1.24
1	A	21	LEU	C-O	-5.10	1.18	1.24
1	A	104	GLU	C-O	-5.08	1.17	1.23
1	A	154	VAL	C-O	-5.06	1.18	1.24
1	A	206	ILE	C-O	-5.04	1.18	1.24
1	A	170	ILE	C-O	-5.04	1.18	1.24
1	B	59	ILE	C-O	-5.03	1.18	1.24
1	A	112	LYS	C-O	-5.02	1.18	1.24
1	A	220	LEU	C-O	-5.02	1.18	1.24

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	217	CYS	N-CA-C	7.20	119.21	111.36
1	B	195	ASN	N-CA-C	6.38	119.22	111.82
1	B	194	SER	N-CA-C	5.86	117.47	111.14
1	B	100	HIS	N-CA-C	5.76	120.46	113.38
1	B	132	ASP	N-CA-C	-5.66	105.11	111.28
1	B	177	THR	CA-C-N	-5.63	112.94	119.47
1	B	177	THR	C-N-CA	-5.63	112.94	119.47
1	A	177	THR	CA-C-N	-5.54	113.00	119.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	177	THR	C-N-CA	-5.54	113.00	119.32
1	A	56	ASP	CA-C-N	-5.53	113.02	119.32
1	A	56	ASP	C-N-CA	-5.53	113.02	119.32
1	A	16	GLY	N-CA-C	5.33	119.02	112.14
1	B	210	GLY	N-CA-C	-5.30	103.40	112.30
1	A	39	VAL	N-CA-C	5.25	115.52	108.17
1	B	74	PHE	N-CA-C	5.21	118.73	112.38
1	B	210	GLY	CA-C-O	-5.00	116.40	122.75

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	1844	0	1849	38	0
1	B	1826	0	1831	70	0
2	A	304	0	0	3	0
2	B	148	0	0	8	0
All	All	4122	0	3680	106	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 (106) 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:14:MET:HE2	1:B:16:GLY:HA2	1.38	1.06
1:B:130:LYS:HA	1:B:167:VAL:HG22	1.44	0.98
1:B:168:TRP:O	1:B:172:THR:HG23	1.74	0.88
1:B:168:TRP:HA	1:B:172:THR:HG21	1.54	0.86
1:A:55:LEU:HD22	1:A:59:ILE:HD12	1.57	0.86
1:B:79:SER:H	1:B:82:MET:HE3	1.41	0.86
1:B:168:TRP:O	1:B:172:THR:CG2	2.23	0.85
1:B:178:PRO:CG	1:B:219:GLU:HB3	2.07	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:LEU:CD2	1:A:59:ILE:HD12	2.07	0.82
1:B:178:PRO:HG2	1:B:219:GLU:CD	2.07	0.80
1:B:178:PRO:HG3	1:B:219:GLU:HB3	1.64	0.78
1:B:164:TYR:CZ	1:B:184:VAL:HG11	2.21	0.76
1:A:45:THR:O	1:A:48:ILE:HG13	1.85	0.76
1:A:193:LYS:HE3	1:A:198:ASP:OD1	1.84	0.76
1:B:182:GLN:NE2	1:B:225:ASP:OD2	2.20	0.74
1:B:130:LYS:HA	1:B:167:VAL:CG2	2.18	0.73
1:B:131:LEU:HB2	1:B:168:TRP:HB3	1.70	0.72
1:B:168:TRP:CA	1:B:172:THR:HG21	2.19	0.72
1:A:45:THR:HA	1:A:48:ILE:HG13	1.73	0.70
1:B:173:GLY:H	1:B:174:LYS:NZ	1.87	0.70
1:B:178:PRO:HG2	1:B:219:GLU:HB3	1.74	0.70
1:B:175:THR:HA	2:B:2119:HOH:O	1.92	0.69
1:B:217:CYS:SG	2:B:2137:HOH:O	2.51	0.68
1:A:14:MET:SD	1:A:21:LEU:HD21	2.35	0.67
1:B:132:ASP:N	1:B:132:ASP:OD1	2.29	0.65
1:B:208:TYR:OH	1:B:210:GLY:HA3	1.97	0.65
1:B:195:ASN:ND2	2:B:2127:HOH:O	2.24	0.64
1:A:25:ILE:HG12	1:A:55:LEU:HG	1.79	0.63
1:B:173:GLY:O	1:B:174:LYS:HG3	1.99	0.63
1:A:55:LEU:HD22	1:A:59:ILE:CD1	2.27	0.62
1:B:178:PRO:HD2	2:B:2139:HOH:O	1.98	0.61
1:A:45:THR:HA	1:A:48:ILE:CG1	2.30	0.61
1:B:164:TYR:CE2	1:B:184:VAL:HG11	2.36	0.61
1:B:208:TYR:CZ	1:B:210:GLY:HA3	2.36	0.61
1:B:54:LYS:NZ	2:B:2060:HOH:O	2.33	0.60
1:B:172:THR:OG1	1:B:174:LYS:NZ	2.30	0.60
1:B:177:THR:HG23	1:B:180:GLN:OE1	2.02	0.59
1:B:173:GLY:N	1:B:174:LYS:NZ	2.50	0.59
1:B:219:GLU:O	1:B:222:SER:OG	2.20	0.58
1:B:218:LYS:NZ	2:B:2138:HOH:O	2.34	0.58
1:B:212:VAL:HG12	2:B:2132:HOH:O	2.03	0.57
1:A:67:TYR:OH	1:A:68:LYS:NZ	2.34	0.56
1:B:96:SER:OG	1:B:170:ILE:HD11	2.06	0.56
1:B:182:GLN:OE1	1:B:223:GLN:HB3	2.05	0.56
1:B:173:GLY:H	1:B:174:LYS:HZ2	1.54	0.56
1:A:164:TYR:CZ	1:A:184:VAL:HG11	2.40	0.55
1:A:56:ASP:OD1	1:A:58:LYS:HB2	2.05	0.55
2:A:2140:HOH:O	1:B:13:LYS:HA	2.07	0.55
1:B:168:TRP:C	1:B:172:THR:CG2	2.79	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:140:GLU:CD	1:B:187:LYS:HD3	2.32	0.54
1:A:25:ILE:HG23	1:A:55:LEU:HA	1.89	0.54
1:A:68:LYS:O	1:A:69:VAL:C	2.51	0.54
1:B:174:LYS:N	1:B:174:LYS:HD2	2.22	0.54
1:B:80:PRO:HB2	1:B:119:GLU:HG3	1.90	0.53
1:A:134:ARG:HD3	1:A:168:TRP:CD2	2.43	0.53
1:A:73:ALA:HA	1:B:97:GLU:CD	2.35	0.52
1:A:237:LYS:HB3	1:A:238:PRO:HD2	1.92	0.52
1:A:45:THR:CA	1:A:48:ILE:HG13	2.38	0.52
1:A:164:TYR:CE2	1:A:184:VAL:HG11	2.46	0.51
1:B:84:LYS:HE2	1:B:119:GLU:O	2.11	0.51
1:A:49:ASP:O	1:A:53:GLN:HG3	2.11	0.51
1:B:178:PRO:HG2	1:B:219:GLU:CB	2.41	0.51
1:A:49:ASP:OD1	1:B:18:LYS:NZ	2.22	0.51
1:B:173:GLY:C	1:B:174:LYS:HD2	2.35	0.51
1:B:173:GLY:N	1:B:174:LYS:HZ3	2.09	0.50
1:B:182:GLN:HE22	1:B:225:ASP:CG	2.19	0.50
1:A:52:ARG:O	1:A:52:ARG:HG3	2.10	0.49
1:B:134:ARG:NH2	2:B:2097:HOH:O	2.44	0.49
1:A:49:ASP:OD1	2:A:2104:HOH:O	2.20	0.49
1:B:168:TRP:HA	1:B:172:THR:CG2	2.34	0.48
1:B:173:GLY:H	1:B:174:LYS:HZ3	1.58	0.48
1:B:178:PRO:HG2	1:B:219:GLU:CG	2.44	0.48
1:B:7:PHE:O	1:B:228:GLY:HA3	2.13	0.48
1:B:58:LYS:HG3	1:B:59:ILE:HG13	1.97	0.47
1:B:12:TRP:HZ2	1:B:28:LEU:HD11	1.80	0.47
1:A:45:THR:C	1:A:48:ILE:HG13	2.38	0.46
1:A:56:ASP:OD1	1:A:58:LYS:N	2.44	0.46
1:A:45:THR:HA	1:A:48:ILE:CD1	2.46	0.46
1:A:45:THR:C	1:A:47:TYR:N	2.73	0.46
1:B:131:LEU:O	1:B:134:ARG:N	2.49	0.46
1:A:40:VAL:HA	1:A:60:ALA:O	2.16	0.45
1:A:182:GLN:CD	1:A:224:PRO:HD2	2.41	0.45
1:B:235:SER:HA	1:B:240:PHE:CD1	2.52	0.45
1:B:173:GLY:C	1:B:174:LYS:CD	2.90	0.45
1:A:208:TYR:CZ	1:A:210:GLY:HA3	2.52	0.45
1:B:67:TYR:HB2	1:B:77:GLU:HB2	1.97	0.45
1:A:186:GLU:OE1	2:A:2251:HOH:O	2.21	0.45
1:B:173:GLY:N	1:B:174:LYS:HZ2	2.14	0.44
1:A:164:TYR:CE1	1:A:184:VAL:HG11	2.52	0.44
1:B:181:ALA:O	1:B:185:HIS:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:LEU:HD12	1:B:206:ILE:HG21	2.00	0.43
1:A:45:THR:HA	1:A:48:ILE:HD11	2.01	0.43
1:B:164:TYR:CE1	1:B:184:VAL:HG11	2.53	0.43
1:A:178:PRO:HG2	1:A:219:GLU:CD	2.43	0.43
1:B:63:ALA:HB3	1:B:83:ILE:HD13	2.01	0.43
1:A:48:ILE:O	1:A:48:ILE:HG22	2.19	0.43
1:B:167:VAL:C	1:B:169:ALA:H	2.27	0.42
1:B:167:VAL:C	1:B:169:ALA:N	2.77	0.42
1:B:68:LYS:O	1:B:69:VAL:C	2.63	0.41
1:A:62:ALA:HB2	1:A:90:TRP:HB2	2.02	0.41
1:B:68:LYS:HG3	1:B:69:VAL:N	2.35	0.41
1:B:168:TRP:C	1:B:172:THR:HG21	2.44	0.41
1:B:12:TRP:O	1:B:14:MET:HG2	2.21	0.41
1:B:84:LYS:HB2	1:B:84:LYS:HE3	1.77	0.40
1:A:16:GLY:O	1:A:17:ARG:HD3	2.21	0.40
1:A:6:PHE:CE1	1:A:228:GLY:HA2	2.56	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	243/250 (97%)	235 (97%)	7 (3%)	1 (0%)	30	27
1	B	240/250 (96%)	234 (98%)	6 (2%)	0	100	100
All	All	483/500 (97%)	469 (97%)	13 (3%)	1 (0%)	43	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	46	ALA

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	193/196 (98%)	188 (97%)	5 (3%)	40	44
1	B	191/196 (97%)	183 (96%)	8 (4%)	26	25
All	All	384/392 (98%)	371 (97%)	13 (3%)	32	33

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	LYS
1	A	71	ASN
1	A	106	ASP
1	A	127	ILE
1	A	202	GLN
1	B	5	LYS
1	B	84	LYS
1	B	107	GLU
1	B	127	ILE
1	B	142	VAL
1	B	167	VAL
1	B	172	THR
1	B	174	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	71	ASN
1	B	100	HIS
1	B	195	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](#)

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	245/250 (98%)	0.06	18 (7%) 21 19	9, 18, 39, 113	0
1	B	242/250 (96%)	0.65	33 (13%) 7 6	13, 31, 84, 99	0
All	All	487/500 (97%)	0.35	51 (10%) 11 10	9, 24, 69, 113	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	48	ILE	9.5
1	B	176	ALA	6.4
1	B	175	THR	5.2
1	A	69	VAL	5.1
1	B	138	ILE	5.0
1	B	137	GLY	4.8
1	B	169	ALA	4.8
1	B	72	GLY	4.3
1	A	46	ALA	4.2
1	B	171	GLY	4.2
1	B	179	GLN	4.1
1	B	69	VAL	4.1
1	B	70	THR	4.1
1	B	139	THR	3.5
1	B	136	ALA	3.5
1	A	72	GLY	3.4
1	B	173	GLY	3.4
1	A	47	TYR	3.4
1	A	70	THR	3.4
1	B	178	PRO	3.4
1	B	67	TYR	3.3
1	B	184	VAL	3.3
1	B	131	LEU	3.3
1	B	168	TRP	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	67	TYR	3.2
1	B	170	ILE	3.2
1	B	181	ALA	3.1
1	A	71	ASN	3.1
1	A	76	GLY	3.0
1	B	76	GLY	3.0
1	B	74	PHE	3.0
1	A	75	THR	2.9
1	A	77	GLU	2.9
1	B	174	LYS	2.9
1	A	53	GLN	2.8
1	A	73	ALA	2.8
1	A	55	LEU	2.8
1	B	177	THR	2.7
1	B	71	ASN	2.7
1	B	172	THR	2.6
1	B	180	GLN	2.5
1	B	68	LYS	2.4
1	B	77	GLU	2.4
1	A	59	ILE	2.4
1	A	51	ALA	2.3
1	B	142	VAL	2.2
1	B	188	LEU	2.2
1	B	219	GLU	2.1
1	A	49	ASP	2.1
1	B	223	GLN	2.0
1	A	32	LYS	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.