



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:09 PM UTC

PDB ID : 2I9E / pdb_00002i9e
Title : Structure of Triosephosphate Isomerase of Tenebrio molitor
Authors : Schmidt, A.; Scheerer, P.; Wessner, H.; Hoehne, W.; Krauss, N.
Deposited on : 2006-09-05
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
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

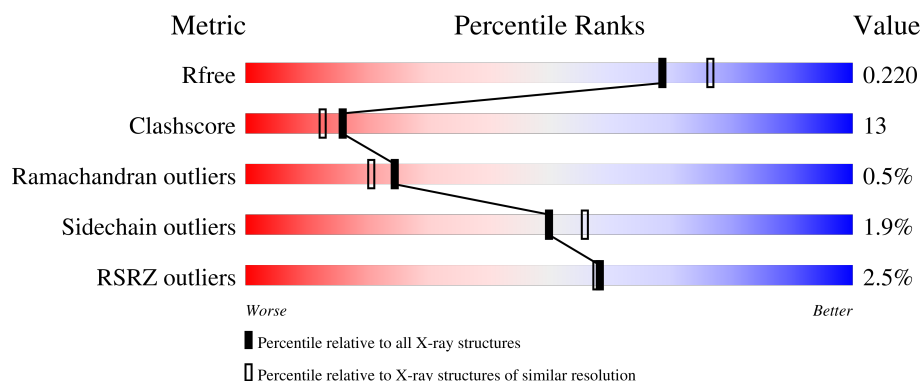
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

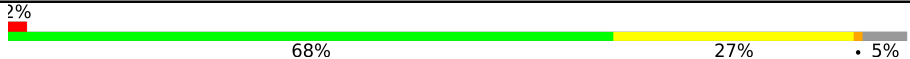
The reported resolution of this entry is 2.00 Å.

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	259	
1	B	259	
1	C	259	
1	D	259	

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	TRS	B	1004	-	-	X	-
2	TRS	D	1003	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7898 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	246	Total	C	N	O	S	0	0	0
			1869	1182	325	354	8			
1	B	248	Total	C	N	O	S	0	0	0
			1884	1193	327	356	8			
1	C	246	Total	C	N	O	S	0	0	0
			1869	1182	325	354	8			
1	D	248	Total	C	N	O	S	0	0	0
			1884	1193	327	356	8			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	241	ASP	ASN	engineered mutation	UNP Q8MPF2
A	243	ILE	VAL	engineered mutation	UNP Q8MPF2
A	246	ARG	LYS	engineered mutation	UNP Q8MPF2
A	248	LEU	-	cloning artifact	UNP Q8MPF2
A	249	VAL	-	cloning artifact	UNP Q8MPF2
A	250	PRO	-	cloning artifact	UNP Q8MPF2
A	251	ARG	-	cloning artifact	UNP Q8MPF2
A	252	GLY	-	cloning artifact	UNP Q8MPF2
A	253	SER	-	cloning artifact	UNP Q8MPF2
A	254	HIS	-	cloning artifact	UNP Q8MPF2
A	255	HIS	-	cloning artifact	UNP Q8MPF2
A	256	HIS	-	cloning artifact	UNP Q8MPF2
A	257	HIS	-	cloning artifact	UNP Q8MPF2
A	258	HIS	-	cloning artifact	UNP Q8MPF2
A	259	HIS	-	cloning artifact	UNP Q8MPF2
B	241	ASP	ASN	engineered mutation	UNP Q8MPF2
B	243	ILE	VAL	engineered mutation	UNP Q8MPF2
B	246	ARG	LYS	engineered mutation	UNP Q8MPF2
B	248	LEU	-	cloning artifact	UNP Q8MPF2
B	249	VAL	-	cloning artifact	UNP Q8MPF2
B	250	PRO	-	cloning artifact	UNP Q8MPF2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	251	ARG	-	cloning artifact	UNP Q8MPF2
B	252	GLY	-	cloning artifact	UNP Q8MPF2
B	253	SER	-	cloning artifact	UNP Q8MPF2
B	254	HIS	-	cloning artifact	UNP Q8MPF2
B	255	HIS	-	cloning artifact	UNP Q8MPF2
B	256	HIS	-	cloning artifact	UNP Q8MPF2
B	257	HIS	-	cloning artifact	UNP Q8MPF2
B	258	HIS	-	cloning artifact	UNP Q8MPF2
B	259	HIS	-	cloning artifact	UNP Q8MPF2
C	241	ASP	ASN	engineered mutation	UNP Q8MPF2
C	243	ILE	VAL	engineered mutation	UNP Q8MPF2
C	246	ARG	LYS	engineered mutation	UNP Q8MPF2
C	248	LEU	-	cloning artifact	UNP Q8MPF2
C	249	VAL	-	cloning artifact	UNP Q8MPF2
C	250	PRO	-	cloning artifact	UNP Q8MPF2
C	251	ARG	-	cloning artifact	UNP Q8MPF2
C	252	GLY	-	cloning artifact	UNP Q8MPF2
C	253	SER	-	cloning artifact	UNP Q8MPF2
C	254	HIS	-	cloning artifact	UNP Q8MPF2
C	255	HIS	-	cloning artifact	UNP Q8MPF2
C	256	HIS	-	cloning artifact	UNP Q8MPF2
C	257	HIS	-	cloning artifact	UNP Q8MPF2
C	258	HIS	-	cloning artifact	UNP Q8MPF2
C	259	HIS	-	cloning artifact	UNP Q8MPF2
D	241	ASP	ASN	engineered mutation	UNP Q8MPF2
D	243	ILE	VAL	engineered mutation	UNP Q8MPF2
D	246	ARG	LYS	engineered mutation	UNP Q8MPF2
D	248	LEU	-	cloning artifact	UNP Q8MPF2
D	249	VAL	-	cloning artifact	UNP Q8MPF2
D	250	PRO	-	cloning artifact	UNP Q8MPF2
D	251	ARG	-	cloning artifact	UNP Q8MPF2
D	252	GLY	-	cloning artifact	UNP Q8MPF2
D	253	SER	-	cloning artifact	UNP Q8MPF2
D	254	HIS	-	cloning artifact	UNP Q8MPF2
D	255	HIS	-	cloning artifact	UNP Q8MPF2
D	256	HIS	-	cloning artifact	UNP Q8MPF2
D	257	HIS	-	cloning artifact	UNP Q8MPF2
D	258	HIS	-	cloning artifact	UNP Q8MPF2
D	259	HIS	-	cloning artifact	UNP Q8MPF2

- Molecule 2 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: C₄H₁₂NO₃).



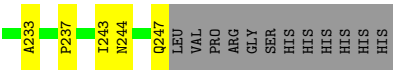
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			8	4	1	3		
2	B	1	Total	C	N	O	0	0
			8	4	1	3		
2	C	1	Total	C	N	O	0	0
			8	4	1	3		
2	D	1	Total	C	N	O	0	0
			8	4	1	3		
2	D	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 3 is water.

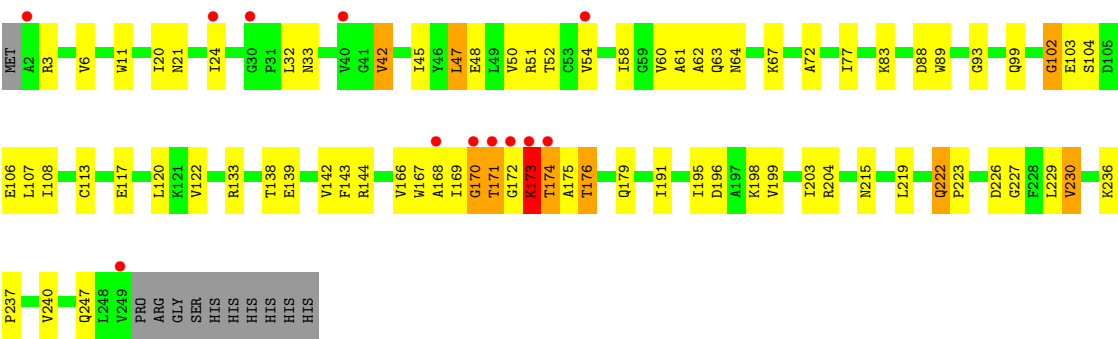
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	93	Total	O	0	0
			93	93		
3	B	87	Total	O	0	0
			87	87		
3	C	95	Total	O	0	0
			95	95		
3	D	77	Total	O	0	0
			77	77		

- Molecule 1: Triosephosphate isomerase





● Molecule 1: Triosephosphate isomerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	37.06Å 145.08Å 94.28Å 90.00° 95.61° 90.00°	Depositor
Resolution (Å)	29.33 – 2.00 29.33 – 2.00	Depositor EDS
% Data completeness (in resolution range)	80.4 (29.33-2.00) 80.4 (29.33-2.00)	Depositor EDS
R_{merge}	0.00	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.219 , 0.239 0.222 , 0.220	Depositor DCC
R_{free} test set	2729 reflections (4.09%)	wwPDB-VP
Wilson B-factor (Å ²)	19.6	Xtriage
Anisotropy	1.032	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7898	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.55% 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: TRS

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	0/1901	0.93	6/2575 (0.2%)
1	B	0.40	0/1916	0.90	5/2596 (0.2%)
1	C	0.38	0/1901	0.93	6/2575 (0.2%)
1	D	0.39	0/1916	0.94	12/2596 (0.5%)
All	All	0.39	0/7634	0.93	29/10342 (0.3%)

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	170	GLY	N-CA-C	-9.71	102.27	112.08
1	A	62	ALA	N-CA-C	-7.98	99.77	110.55
1	C	62	ALA	N-CA-C	-7.75	100.08	110.55
1	B	122	VAL	N-CA-C	7.07	118.99	108.46
1	D	122	VAL	N-CA-C	6.96	118.83	108.46
1	D	62	ALA	N-CA-C	-6.62	100.72	110.52
1	B	62	ALA	N-CA-C	-6.57	100.80	110.52
1	B	229	LEU	N-CA-C	-6.56	94.69	107.62
1	C	102	GLY	N-CA-C	6.54	123.03	113.48
1	D	203	ILE	N-CA-C	6.45	117.40	108.89
1	C	42	VAL	N-CA-C	6.25	114.78	107.84
1	D	42	VAL	N-CA-C	6.07	114.58	107.84
1	A	102	GLY	N-CA-C	6.07	122.34	113.48
1	A	135	ALA	N-CA-C	-6.05	105.76	113.02
1	C	135	ALA	N-CA-C	-5.96	106.05	113.15
1	D	176	THR	N-CA-C	-5.93	103.31	110.13
1	A	42	VAL	N-CA-C	5.84	114.32	107.84
1	B	42	VAL	N-CA-C	5.82	114.30	107.84
1	D	67	LYS	N-CA-C	5.72	119.96	112.92
1	D	222	GLN	CA-C-N	5.58	125.68	119.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	222	GLN	C-N-CA	5.58	125.68	119.32
1	D	230	VAL	N-CA-C	5.57	116.79	109.21
1	B	67	LYS	N-CA-C	5.56	119.76	112.92
1	A	122	VAL	N-CA-C	5.52	117.36	108.85
1	A	67	LYS	N-CA-C	5.47	120.07	113.17
1	C	67	LYS	N-CA-C	5.37	119.94	113.17
1	C	122	VAL	N-CA-C	5.36	117.10	108.85
1	D	229	LEU	N-CA-C	-5.27	97.09	107.69
1	D	102	GLY	N-CA-C	5.20	121.07	113.48

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	1869	0	1885	44	0
1	B	1884	0	1905	54	0
1	C	1869	0	1885	48	0
1	D	1884	0	1905	59	0
2	B	16	0	24	6	0
2	C	8	0	12	4	0
2	D	16	0	24	6	0
3	A	93	0	0	1	0
3	B	87	0	0	2	0
3	C	95	0	0	1	0
3	D	77	0	0	0	0
All	All	7898	0	7640	201	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 (201) 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:99:GLN:HE22	1:B:166:VAL:HG11	1.17	1.04
1:B:232:GLY:HA2	2:B:1004:TRS:H22	1.51	0.93
1:D:99:GLN:HE22	1:D:166:VAL:HG11	1.34	0.91
1:B:131:GLU:HG2	3:B:1012:HOH:O	1.80	0.79
1:C:102:GLY:HA3	2:C:1001:TRS:H32	1.66	0.78
1:C:171:THR:HG22	1:C:173:LYS:H	1.49	0.78
1:A:171:THR:HG22	1:A:173:LYS:H	1.48	0.76
1:D:20:ILE:O	1:D:24:ILE:HG12	1.86	0.74
1:B:20:ILE:O	1:B:24:ILE:HG12	1.86	0.74
1:C:20:ILE:O	1:C:24:ILE:HG12	1.90	0.71
1:D:103:GLU:HA	2:D:1003:TRS:H32	1.72	0.71
1:D:215:ASN:O	1:D:219:LEU:HD13	1.90	0.71
1:B:99:GLN:NE2	1:B:166:VAL:HG11	1.99	0.70
1:A:20:ILE:O	1:A:24:ILE:HG12	1.90	0.69
1:C:168:ALA:O	1:C:171:THR:HB	1.93	0.68
1:A:168:ALA:O	1:A:171:THR:HB	1.93	0.67
1:D:176:THR:OG1	1:D:179:GLN:HG3	1.97	0.65
1:D:107:LEU:CD2	2:D:1003:TRS:N	2.60	0.64
1:B:196:ASP:OD2	1:B:199:VAL:HG23	2.01	0.61
1:B:51:ARG:HA	1:B:60:VAL:HG21	1.83	0.60
1:D:51:ARG:HA	1:D:60:VAL:HG21	1.83	0.59
1:D:191:ILE:HG23	1:D:195:ILE:HD12	1.85	0.59
1:C:93:GLY:HA2	1:C:108:ILE:HD12	1.85	0.59
1:A:6:VAL:HG22	1:A:204:ARG:NH1	2.18	0.59
1:B:99:GLN:HE22	1:B:166:VAL:CG1	2.04	0.59
1:B:177:PRO:HG3	1:B:219:LEU:HD12	1.84	0.59
1:C:6:VAL:HG22	1:C:204:ARG:NH1	2.18	0.58
1:B:231:GLY:C	2:B:1004:TRS:H12	2.29	0.58
1:D:173:LYS:HD3	1:D:174:THR:N	2.17	0.58
1:A:93:GLY:HA2	1:A:108:ILE:HD12	1.86	0.57
1:D:107:LEU:HD22	2:D:1003:TRS:HN3	1.70	0.57
1:B:26:PHE:HD1	3:B:1083:HOH:O	1.88	0.56
1:B:171:THR:CG2	1:B:173:LYS:O	2.53	0.56
1:C:151:ALA:C	1:C:152:LYS:HD3	2.31	0.56
1:B:33:ASN:HD22	1:B:247:GLN:CG	2.19	0.55
1:C:99:GLN:HE22	1:C:166:VAL:HG11	1.71	0.55
1:A:33:ASN:HB2	1:A:244:ASN:OD1	2.05	0.55
1:B:177:PRO:HG3	1:B:219:LEU:CD1	2.36	0.55
1:B:171:THR:HG23	1:B:173:LYS:H	1.72	0.55
1:B:232:GLY:N	2:B:1004:TRS:H12	2.22	0.55
1:A:99:GLN:HE22	1:A:166:VAL:HG11	1.71	0.54
1:A:151:ALA:C	1:A:152:LYS:HD3	2.31	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:33:ASN:HB2	1:C:244:ASN:OD1	2.07	0.54
1:A:6:VAL:CG2	1:A:204:ARG:HG2	2.38	0.54
1:B:6:VAL:O	1:B:227:GLY:HA3	2.08	0.53
1:C:6:VAL:CG2	1:C:204:ARG:HG2	2.37	0.53
1:A:176:THR:OG1	1:A:179:GLN:HG3	2.09	0.53
1:B:6:VAL:CG2	1:B:204:ARG:NH1	2.71	0.53
1:D:168:ALA:O	1:D:169:ILE:C	2.52	0.53
1:D:107:LEU:CD2	2:D:1003:TRS:HN1	2.21	0.52
1:B:48:GLU:O	1:B:52:THR:HG23	2.10	0.52
1:D:42:VAL:HG11	1:D:50:VAL:HG21	1.91	0.52
1:D:6:VAL:O	1:D:227:GLY:HA3	2.10	0.52
1:D:83:LYS:HG2	1:D:120:LEU:HD21	1.92	0.52
1:D:173:LYS:HD3	1:D:173:LYS:C	2.35	0.52
1:D:48:GLU:O	1:D:52:THR:HG23	2.10	0.52
1:B:42:VAL:HG11	1:B:50:VAL:HG21	1.92	0.51
1:A:104:SER:O	1:A:108:ILE:HG12	2.11	0.51
1:B:104:SER:O	1:B:108:ILE:HG12	2.10	0.51
1:D:174:THR:HG22	1:D:175:ALA:N	2.26	0.50
1:B:232:GLY:CA	2:B:1004:TRS:H22	2.32	0.50
1:A:51:ARG:HA	1:A:60:VAL:HG21	1.93	0.50
1:C:26:PHE:CE2	1:C:237:PRO:HG3	2.46	0.50
1:D:54:VAL:CG1	1:D:58:ILE:HB	2.42	0.50
1:A:45:ILE:HD11	1:B:77:ILE:HD12	1.92	0.50
1:C:104:SER:O	1:C:108:ILE:HG12	2.12	0.50
1:C:176:THR:OG1	1:C:179:GLN:HG3	2.10	0.50
1:D:196:ASP:OD2	1:D:198:LYS:HB2	2.10	0.50
1:D:33:ASN:HD22	1:D:247:GLN:HG2	1.77	0.50
1:D:93:GLY:HA2	1:D:108:ILE:HD12	1.94	0.50
1:B:54:VAL:CG1	1:B:58:ILE:HB	2.42	0.50
1:C:191:ILE:HG23	1:C:195:ILE:HD12	1.92	0.49
1:D:196:ASP:HB3	1:D:199:VAL:HG23	1.93	0.49
1:A:26:PHE:CE2	1:A:237:PRO:HG3	2.47	0.49
1:C:51:ARG:HA	1:C:60:VAL:HG21	1.92	0.49
1:A:191:ILE:HG23	1:A:195:ILE:HD12	1.93	0.49
1:D:167:TRP:CZ2	1:D:175:ALA:HB2	2.46	0.49
1:D:107:LEU:HD22	2:D:1003:TRS:N	2.25	0.49
1:A:12:LYS:HD3	1:B:72:ALA:HA	1.95	0.49
1:B:138:THR:O	1:B:142:VAL:HG22	2.13	0.49
1:C:113:CYS:O	1:C:117:GLU:HG3	2.13	0.49
1:D:107:LEU:CD2	2:D:1003:TRS:HN3	2.22	0.48
1:A:113:CYS:O	1:A:117:GLU:HG3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:102:GLY:CA	2:C:1001:TRS:H32	2.42	0.48
1:D:138:THR:O	1:D:142:VAL:HG22	2.14	0.47
1:C:26:PHE:CZ	1:C:237:PRO:HG3	2.49	0.47
1:C:6:VAL:HG23	1:C:204:ARG:HG2	1.95	0.47
1:B:192:CYS:SG	1:B:200:GLY:HA3	2.55	0.47
1:B:215:ASN:O	1:B:219:LEU:HD13	2.14	0.47
1:A:26:PHE:CZ	1:A:237:PRO:HG3	2.50	0.47
1:C:45:ILE:HD11	1:D:77:ILE:HD12	1.97	0.47
1:D:3:ARG:HE	1:D:226:ASP:CG	2.23	0.47
1:D:113:CYS:O	1:D:117:GLU:HG3	2.14	0.47
1:A:6:VAL:CG2	1:A:204:ARG:NH1	2.78	0.47
1:A:6:VAL:O	1:A:227:GLY:HA3	2.16	0.46
1:A:6:VAL:HG23	1:A:204:ARG:HG2	1.95	0.46
1:C:6:VAL:CG2	1:C:204:ARG:NH1	2.79	0.46
1:B:168:ALA:O	1:B:171:THR:HG22	2.15	0.46
1:D:104:SER:O	1:D:108:ILE:HG12	2.15	0.46
1:B:24:ILE:HD12	1:B:54:VAL:HG22	1.96	0.46
1:A:63:GLN:O	1:A:64:ASN:HB2	2.16	0.46
2:C:1001:TRS:C2	1:D:102:GLY:HA3	2.46	0.46
1:A:108:ILE:O	1:A:112:VAL:HG23	2.15	0.46
1:C:108:ILE:O	1:C:112:VAL:HG23	2.16	0.46
1:C:63:GLN:O	1:C:64:ASN:HB2	2.15	0.46
1:D:61:ALA:HB2	1:D:89:TRP:HB2	1.98	0.45
1:A:11:TRP:CZ3	1:A:20:ILE:HG12	2.51	0.45
1:C:11:TRP:CZ3	1:C:20:ILE:HG12	2.51	0.45
1:C:61:ALA:HB2	1:C:89:TRP:CB	2.46	0.45
1:D:24:ILE:HD12	1:D:54:VAL:HG22	1.97	0.45
1:D:133:ARG:HD2	1:D:167:TRP:CG	2.52	0.45
1:A:36:THR:HG21	1:A:243:ILE:HG23	1.99	0.45
1:B:133:ARG:HD2	1:B:167:TRP:CG	2.51	0.45
1:C:6:VAL:O	1:C:227:GLY:HA3	2.15	0.45
1:A:141:VAL:O	1:A:145:GLN:HG3	2.16	0.45
1:B:113:CYS:O	1:B:117:GLU:HG3	2.16	0.45
1:C:61:ALA:HB2	1:C:89:TRP:HB2	1.97	0.45
1:A:48:GLU:O	1:A:52:THR:HG23	2.17	0.45
1:D:139:GLU:HB3	1:D:143:PHE:CE2	2.52	0.45
1:D:196:ASP:HB3	1:D:199:VAL:CG2	2.47	0.45
1:B:61:ALA:HB2	1:B:89:TRP:HB2	1.98	0.45
1:B:88:ASP:OD1	1:B:88:ASP:N	2.50	0.45
1:B:139:GLU:HB3	1:B:143:PHE:CE2	2.52	0.45
2:C:1001:TRS:H22	1:D:102:GLY:HA3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:12:LYS:HD3	1:D:72:ALA:HA	1.98	0.45
1:D:169:ILE:HD12	1:D:169:ILE:N	2.32	0.45
1:D:172:GLY:O	1:D:173:LYS:C	2.60	0.45
1:B:209:GLY:O	1:B:211:VAL:HG23	2.17	0.44
1:C:36:THR:HG21	1:C:243:ILE:HG23	1.99	0.44
1:D:169:ILE:H	1:D:169:ILE:CD1	2.31	0.44
1:C:48:GLU:O	1:C:52:THR:HG23	2.17	0.44
1:D:106:GLU:CD	1:D:106:GLU:N	2.76	0.44
1:B:12:LYS:HD2	2:B:1004:TRS:H32	1.99	0.44
1:C:82:ILE:CG2	1:C:87:ALA:HB3	2.48	0.44
1:A:82:ILE:CG2	1:A:87:ALA:HB3	2.47	0.44
1:B:24:ILE:HD12	1:B:54:VAL:CG2	2.48	0.44
1:C:230:VAL:CG1	1:C:233:ALA:HB3	2.48	0.44
1:A:196:ASP:OD2	1:A:198:LYS:HB2	2.18	0.44
1:D:32:LEU:HD23	1:D:240:VAL:HG13	2.00	0.44
1:D:144:ARG:HH21	1:D:144:ARG:HG2	1.83	0.44
1:C:126:ILE:HB	1:C:145:GLN:OE1	2.18	0.43
1:A:230:VAL:CG1	1:A:233:ALA:HB3	2.48	0.43
1:C:26:PHE:HD1	3:C:1080:HOH:O	2.02	0.43
1:D:33:ASN:HD22	1:D:247:GLN:CG	2.32	0.43
1:B:232:GLY:HA2	2:B:1004:TRS:C2	2.36	0.43
1:D:24:ILE:HD12	1:D:54:VAL:CG2	2.48	0.43
1:B:191:ILE:HG23	1:B:195:ILE:HD12	2.00	0.43
1:C:129:THR:OG1	1:C:132:GLU:HG3	2.18	0.43
1:C:196:ASP:OD2	1:C:198:LYS:HB2	2.18	0.43
1:D:47:LEU:HD23	1:D:47:LEU:HA	1.79	0.43
1:C:54:VAL:CG1	1:C:58:ILE:HB	2.49	0.43
1:D:170:GLY:C	1:D:172:GLY:N	2.74	0.43
1:B:215:ASN:C	1:B:215:ASN:OD1	2.62	0.43
1:A:61:ALA:HB2	1:A:89:TRP:CB	2.50	0.42
1:C:163:TYR:CZ	1:C:183:VAL:HG11	2.54	0.42
1:D:88:ASP:OD1	1:D:88:ASP:N	2.52	0.42
1:A:25:GLY:O	1:A:29:SER:HB3	2.20	0.42
1:A:61:ALA:HB2	1:A:89:TRP:HB2	2.01	0.42
1:B:26:PHE:CE2	1:B:237:PRO:HG3	2.54	0.42
1:C:5:PHE:CE1	1:C:227:GLY:HA2	2.55	0.42
1:C:150:ALA:HA	1:C:156:TRP:CZ2	2.54	0.42
1:D:6:VAL:CG2	1:D:204:ARG:NH1	2.82	0.42
1:D:171:THR:O	1:D:173:LYS:N	2.53	0.42
1:B:33:ASN:HB2	1:B:247:GLN:OE1	2.19	0.42
1:B:89:TRP:CD2	1:B:121:LYS:HB2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:141:VAL:O	1:C:145:GLN:HG3	2.19	0.42
1:A:54:VAL:CG1	1:A:58:ILE:HB	2.49	0.42
1:C:163:TYR:CE1	1:C:183:VAL:HG11	2.55	0.42
1:D:169:ILE:HD12	1:D:169:ILE:H	1.84	0.42
1:A:163:TYR:CE1	1:A:183:VAL:HG11	2.55	0.42
1:A:150:ALA:HA	1:A:156:TRP:CZ2	2.55	0.41
1:B:83:LYS:HG2	1:B:120:LEU:HD21	2.01	0.41
1:C:106:GLU:H	1:C:106:GLU:CD	2.28	0.41
1:A:106:GLU:H	1:A:106:GLU:CD	2.28	0.41
1:A:126:ILE:HB	1:A:145:GLN:OE1	2.21	0.41
1:A:129:THR:OG1	1:A:132:GLU:HG3	2.20	0.41
1:B:11:TRP:CE3	1:B:42:VAL:HG12	2.55	0.41
1:B:176:THR:OG1	1:B:179:GLN:HG3	2.20	0.41
1:D:63:GLN:O	1:D:64:ASN:HB2	2.21	0.41
1:C:161:ILE:HG13	1:C:203:ILE:HD11	2.02	0.41
1:D:236:LYS:HB3	1:D:237:PRO:CD	2.50	0.41
1:B:144:ARG:HG2	1:B:144:ARG:HH21	1.84	0.41
1:D:11:TRP:CE3	1:D:42:VAL:HG12	2.56	0.41
1:A:88:ASP:OD1	1:A:88:ASP:N	2.54	0.41
1:B:63:GLN:O	1:B:64:ASN:HB2	2.20	0.41
1:D:198:LYS:HE2	1:D:198:LYS:HB3	1.83	0.41
1:A:244:ASN:HB3	1:A:247:GLN:OE1	2.21	0.41
1:C:229:LEU:HD12	1:C:229:LEU:HA	1.90	0.41
1:D:236:LYS:HB3	1:D:237:PRO:HD2	2.03	0.41
1:B:6:VAL:HG22	1:B:204:ARG:NH1	2.36	0.40
1:A:163:TYR:CZ	1:A:183:VAL:HG11	2.56	0.40
1:B:33:ASN:OD1	1:B:35:ASP:N	2.54	0.40
1:A:106:GLU:O	1:A:110:GLU:HG3	2.20	0.40
1:A:246:ARG:HB2	3:A:327:HOH:O	2.21	0.40
1:B:6:VAL:HG23	1:B:204:ARG:HG2	2.04	0.40
1:B:47:LEU:HA	1:B:47:LEU:HD23	1.80	0.40
1:C:244:ASN:HB3	1:C:247:GLN:OE1	2.21	0.40
1:B:141:VAL:O	1:B:145:GLN:HG3	2.21	0.40
1:C:88:ASP:OD1	1:C:88:ASP:N	2.54	0.40
1:C:150:ALA:HA	1:C:156:TRP:HZ2	1.86	0.40
1:D:222:GLN:HA	1:D:223:PRO:HD3	1.84	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	244/259 (94%)	236 (97%)	7 (3%)	1 (0%)	30	27
1	B	246/259 (95%)	236 (96%)	10 (4%)	0	100	100
1	C	244/259 (94%)	236 (97%)	6 (2%)	2 (1%)	16	11
1	D	246/259 (95%)	234 (95%)	10 (4%)	2 (1%)	16	11
All	All	980/1036 (95%)	942 (96%)	33 (3%)	5 (0%)	24	21

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	173	LYS
1	D	174	THR
1	A	30	GLY
1	C	29	SER
1	C	30	GLY

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	197/209 (94%)	194 (98%)	3 (2%)	57	64
1	B	199/209 (95%)	196 (98%)	3 (2%)	57	64
1	C	197/209 (94%)	194 (98%)	3 (2%)	57	64
1	D	199/209 (95%)	193 (97%)	6 (3%)	36	38
All	All	792/836 (95%)	777 (98%)	15 (2%)	50	56

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	VAL
1	A	171	THR
1	A	219	LEU
1	B	21	ASN
1	B	45	ILE
1	B	47	LEU
1	C	6	VAL
1	C	171	THR
1	C	219	LEU
1	D	21	ASN
1	D	45	ILE
1	D	47	LEU
1	D	171	THR
1	D	173	LYS
1	D	230	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	99	GLN
1	A	114	HIS
1	A	158	ASN
1	A	201	ASN
1	B	14	ASN
1	B	99	GLN
1	B	184	HIS
1	B	206	GLN
1	C	21	ASN
1	C	99	GLN
1	C	114	HIS
1	C	158	ASN
1	C	184	HIS
1	C	201	ASN
1	D	99	GLN
1	D	184	HIS
1	D	194	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](#)

5 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	TRS	C	1001	-	7,7,7	0.52	0	9,9,9	0.49	0
2	TRS	B	1002	-	7,7,7	0.62	0	9,9,9	0.46	0
2	TRS	B	1004	-	7,7,7	0.67	0	9,9,9	0.66	0
2	TRS	D	1005	-	7,7,7	0.59	0	9,9,9	0.56	0
2	TRS	D	1003	-	7,7,7	0.58	0	9,9,9	0.57	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	TRS	C	1001	-	-	0/9/9/9	-
2	TRS	B	1002	-	-	0/9/9/9	-
2	TRS	B	1004	-	-	0/9/9/9	-
2	TRS	D	1005	-	-	0/9/9/9	-
2	TRS	D	1003	-	-	0/9/9/9	-

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.

3 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1001	TRS	4	0
2	B	1004	TRS	6	0
2	D	1003	TRS	6	0

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	246/259 (94%)	0.08	5 (2%) 65 64	16, 26, 46, 61	0
1	B	248/259 (95%)	0.09	6 (2%) 59 59	17, 26, 45, 58	0
1	C	246/259 (94%)	0.12	2 (0%) 82 82	17, 28, 47, 58	0
1	D	248/259 (95%)	0.23	12 (4%) 35 34	18, 27, 48, 71	0
All	All	988/1036 (95%)	0.13	25 (2%) 58 58	16, 27, 47, 71	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	171	THR	5.9
1	D	174	THR	4.4
1	D	172	GLY	4.0
1	B	30	GLY	4.0
1	B	249	VAL	3.9
1	D	170	GLY	3.7
1	B	2	ALA	3.5
1	A	31	PRO	3.5
1	D	54	VAL	3.4
1	D	249	VAL	3.4
1	D	173	LYS	3.2
1	D	30	GLY	3.1
1	C	30	GLY	3.0
1	C	29	SER	3.0
1	A	32	LEU	2.9
1	B	26	PHE	2.8
1	A	54	VAL	2.4
1	D	24	ILE	2.4
1	D	40	VAL	2.3
1	A	30	GLY	2.3
1	D	168	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	170	GLY	2.2
1	B	169	ILE	2.1
1	D	2	ALA	2.1
1	A	26	PHE	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	TRS	D	1005	8/8	0.46	0.29	68,79,84,84	0
2	TRS	D	1003	8/8	0.69	0.22	56,61,66,67	0
2	TRS	B	1004	8/8	0.76	0.18	23,37,45,52	0
2	TRS	B	1002	8/8	0.79	0.16	48,54,59,60	0
2	TRS	C	1001	8/8	0.79	0.19	53,68,73,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.