



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 10:59 pm BST

PDB ID : 3TIM
Title : THE CRYSTAL STRUCTURE OF THE "OPEN" AND THE "CLOSED"
CONFORMATION OF THE FLEXIBLE LOOP OF TRYPANOSOMAL
TRIOSEPHOSPHATE ISOMERASE
Authors : Wierenga, R.K.
Deposited on : 1990-05-15
Resolution : 2.80 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 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	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

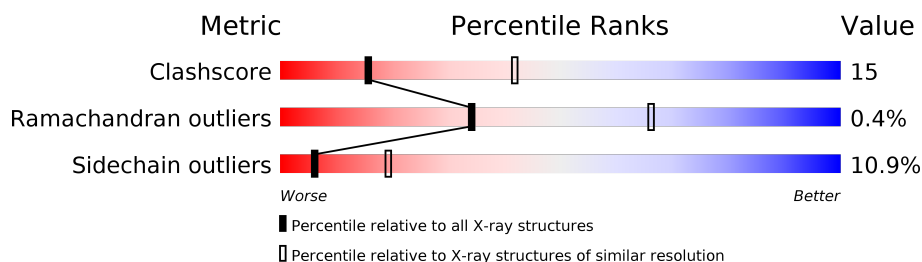
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)

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 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	250	
1	B	250	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3808 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	249	Total	C	N	O	S	0	0	0
			1889	1200	334	350	5			
1	B	249	Total	C	N	O	S	0	0	0
			1889	1200	334	350	5			

- Molecule 2 is water.

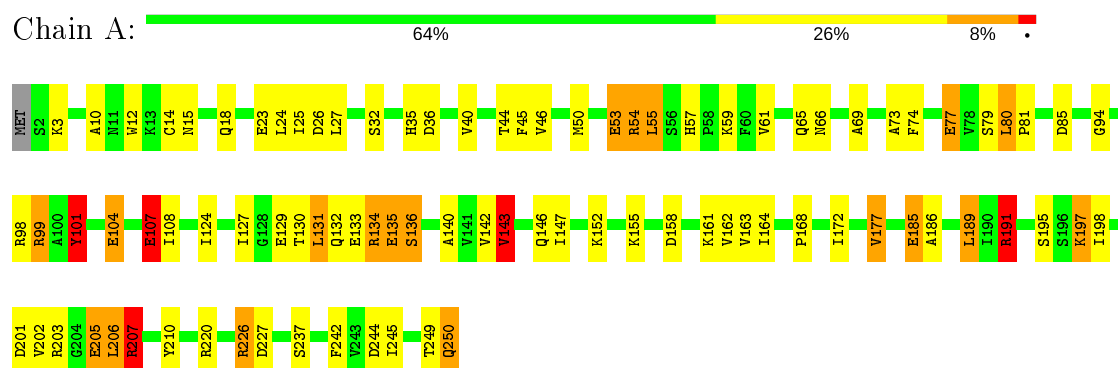
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	23	Total	O	0	0
			23	23		
2	B	7	Total	O	0	0
			7	7		

3 Residue-property plots

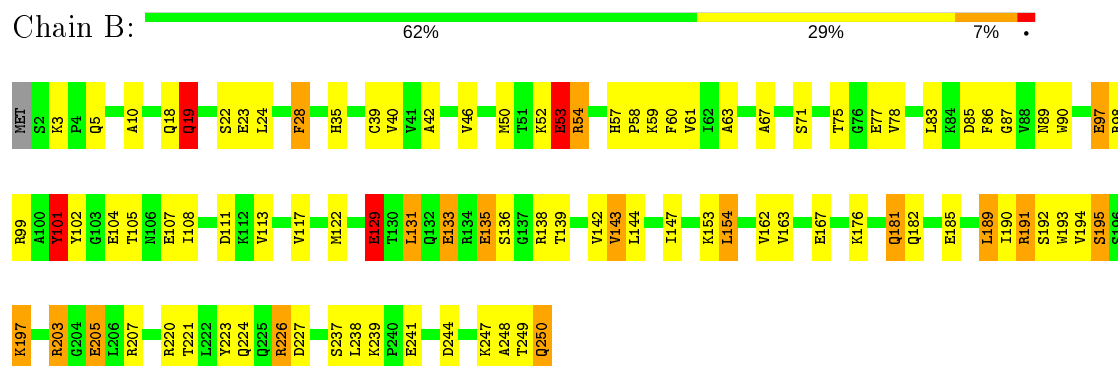
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according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



• Molecule 1: TRIOSEPHOSPHATE ISOMERASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	113.14Å 96.48Å 46.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.80	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-2.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.139 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3808	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

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	0.99	9/1923 (0.5%)	1.52	27/2606 (1.0%)
1	B	1.00	13/1923 (0.7%)	1.40	19/2606 (0.7%)
All	All	1.00	22/3846 (0.6%)	1.46	46/5212 (0.9%)

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	1	1
All	All	1	3

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	133	GLU	CD-OE2	8.83	1.35	1.25
1	A	107	GLU	CD-OE2	8.03	1.34	1.25
1	B	167	GLU	CD-OE2	7.83	1.34	1.25
1	B	205	GLU	CD-OE2	7.83	1.34	1.25
1	A	53	GLU	CD-OE2	7.45	1.33	1.25
1	B	53	GLU	CD-OE2	7.44	1.33	1.25
1	B	185	GLU	CD-OE2	7.18	1.33	1.25
1	B	129	GLU	CD-OE2	7.17	1.33	1.25
1	A	129	GLU	CD-OE2	7.05	1.33	1.25
1	A	104	GLU	CD-OE2	7.01	1.33	1.25
1	A	135	GLU	CD-OE2	6.75	1.33	1.25
1	B	104	GLU	CD-OE2	6.56	1.32	1.25
1	A	23	GLU	CD-OE2	6.54	1.32	1.25
1	B	107	GLU	CD-OE2	6.43	1.32	1.25
1	B	135	GLU	CD-OE2	6.41	1.32	1.25
1	A	205	GLU	CD-OE2	6.40	1.32	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	185	GLU	CD-OE2	6.26	1.32	1.25
1	B	241	GLU	CD-OE2	5.78	1.32	1.25
1	B	97	GLU	CD-OE2	5.74	1.31	1.25
1	B	77	GLU	CD-OE2	5.49	1.31	1.25
1	A	77	GLU	CD-OE2	5.46	1.31	1.25
1	B	23	GLU	CD-OE2	5.41	1.31	1.25

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	134	ARG	NE-CZ-NH2	-11.52	114.54	120.30
1	A	191	ARG	NE-CZ-NH1	9.55	125.07	120.30
1	A	134	ARG	NE-CZ-NH1	9.38	124.99	120.30
1	A	207	ARG	NE-CZ-NH1	-8.85	115.88	120.30
1	B	99	ARG	NE-CZ-NH2	-8.37	116.11	120.30
1	A	226	ARG	NE-CZ-NH2	-8.20	116.20	120.30
1	B	101	TYR	CB-CG-CD2	-7.90	116.26	121.00
1	A	98	ARG	NE-CZ-NH1	7.40	124.00	120.30
1	A	226	ARG	NE-CZ-NH1	7.32	123.96	120.30
1	A	36	ASP	CB-CG-OD1	7.06	124.65	118.30
1	B	85	ASP	CB-CG-OD2	-6.99	112.01	118.30
1	A	244	ASP	CB-CG-OD2	-6.94	112.06	118.30
1	A	36	ASP	CB-CG-OD2	-6.81	112.17	118.30
1	A	189	LEU	CB-CA-C	-6.56	97.73	110.20
1	A	210	TYR	CB-CG-CD2	-6.56	117.07	121.00
1	A	163	VAL	CA-CB-CG2	-6.35	101.37	110.90
1	B	226	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	A	227	ASP	CB-CG-OD1	6.16	123.84	118.30
1	A	244	ASP	CB-CG-OD1	6.12	123.81	118.30
1	B	244	ASP	CB-CG-OD2	-5.98	112.92	118.30
1	B	85	ASP	CB-CG-OD1	5.90	123.61	118.30
1	B	220	ARG	NE-CZ-NH1	5.84	123.22	120.30
1	A	227	ASP	CB-CG-OD2	-5.84	113.05	118.30
1	A	201	ASP	CB-CG-OD1	5.81	123.53	118.30
1	A	26	ASP	CB-CG-OD2	-5.73	113.14	118.30
1	A	85	ASP	CB-CG-OD2	-5.65	113.21	118.30
1	B	58	PRO	N-CA-CB	5.64	110.06	103.30
1	A	101	TYR	CB-CG-CD2	-5.63	117.62	121.00
1	A	177	VAL	CB-CA-C	-5.62	100.73	111.40
1	B	249	THR	N-CA-CB	5.61	120.96	110.30
1	A	143	VAL	N-CA-CB	5.60	123.82	111.50
1	B	197	LYS	N-CA-C	5.60	126.13	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	85	ASP	CB-CG-OD1	5.58	123.32	118.30
1	B	101	TYR	CA-CB-CG	5.57	123.98	113.40
1	B	250	GLN	CB-CA-C	-5.47	99.46	110.40
1	A	140	ALA	N-CA-CB	5.43	117.70	110.10
1	B	227	ASP	CB-CG-OD2	-5.39	113.45	118.30
1	B	220	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	A	101	TYR	CB-CG-CD1	5.35	124.21	121.00
1	B	143	VAL	CG1-CB-CG2	-5.32	102.39	110.90
1	B	111	ASP	CB-CG-OD2	-5.30	113.53	118.30
1	B	54	ARG	NE-CZ-NH1	-5.29	117.66	120.30
1	A	207	ARG	NE-CZ-NH2	5.28	122.94	120.30
1	B	129	GLU	N-CA-CB	5.19	119.94	110.60
1	A	158	ASP	CB-CG-OD1	5.17	122.95	118.30
1	B	28	PHE	N-CA-CB	5.10	119.78	110.60

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	249	THR	CA

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	101	TYR	Sidechain
1	A	124	ILE	Mainchain
1	B	101	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	1889	0	1925	59	5
1	B	1889	0	1925	60	0
2	A	23	0	0	1	0
2	B	7	0	0	0	0
All	All	3808	0	3850	112	5

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 15.

All (112) 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:57:HIS:HD2	1:B:59:LYS:H	1.20	0.85
1:A:191:ARG:HD3	1:A:203:ARG:HD2	1.62	0.80
1:B:52:LYS:NZ	1:B:86:PHE:O	2.15	0.76
1:A:143:VAL:O	1:A:147:ILE:HG22	1.87	0.74
1:B:52:LYS:HZ1	1:B:87:GLY:HA3	1.52	0.74
1:B:5:GLN:HE21	1:B:207:ARG:HH12	1.36	0.72
1:A:79:SER:OG	1:A:81:PRO:HD2	1.90	0.72
1:A:99:ARG:HH21	1:A:99:ARG:HG2	1.56	0.69
1:A:195:SER:HB2	1:A:203:ARG:HD3	1.74	0.69
1:B:5:GLN:HE21	1:B:207:ARG:NH1	1.91	0.68
1:B:105:THR:OG1	1:B:108:ILE:HG12	1.93	0.68
1:A:73:ALA:HB1	1:B:97:GLU:OE2	1.94	0.68
1:B:57:HIS:CD2	1:B:59:LYS:H	2.08	0.67
1:A:50:MET:O	1:A:54:ARG:HB2	1.95	0.67
1:B:5:GLN:O	1:B:207:ARG:HD2	1.96	0.65
1:A:57:HIS:HD2	1:A:59:LYS:H	1.43	0.65
1:B:190:ILE:O	1:B:194:VAL:HG23	1.98	0.64
1:B:5:GLN:NE2	1:B:207:ARG:HH12	1.95	0.64
1:B:139:THR:O	1:B:143:VAL:HG22	1.99	0.63
1:A:10:ALA:HB1	1:A:237:SER:HB2	1.82	0.61
1:B:226:ARG:HG3	1:B:226:ARG:O	2.00	0.61
1:A:130:THR:HG22	1:A:133:GLU:OE2	2.01	0.61
1:B:113:VAL:O	1:B:117:VAL:HG23	2.01	0.60
1:B:3:LYS:HG2	1:B:226:ARG:HH21	1.65	0.60
1:B:248:ALA:O	1:B:250:GLN:HG2	2.01	0.60
1:A:131:LEU:HD22	1:A:135:GLU:HG2	1.84	0.59
1:A:164:ILE:HD12	1:A:206:LEU:HD11	1.84	0.59
1:B:40:VAL:HG22	1:B:61:VAL:HG22	1.83	0.58
1:A:107:GLU:HG2	1:A:108:ILE:N	2.18	0.58
1:B:52:LYS:NZ	1:B:87:GLY:HA3	2.18	0.58
1:A:27:LEU:O	1:A:27:LEU:HD23	2.04	0.58
1:B:162:VAL:HG12	1:B:163:VAL:N	2.19	0.57
1:A:57:HIS:CD2	1:A:59:LYS:HB2	2.40	0.57
1:A:132:GLN:O	1:A:136:SER:OG	2.23	0.56
1:A:191:ARG:HD3	1:A:203:ARG:CD	2.35	0.56
1:A:79:SER:CB	1:A:81:PRO:HD2	2.35	0.56
1:B:133:GLU:HG2	1:B:138:ARG:NH1	2.21	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:LYS:NZ	1:B:223:TYR:O	2.39	0.56
1:B:52:LYS:CE	1:B:86:PHE:O	2.54	0.56
1:A:57:HIS:CD2	1:A:59:LYS:H	2.22	0.55
1:A:250:GLN:NE2	1:A:250:GLN:HA	2.12	0.55
1:A:77:GLU:OE2	1:B:98:ARG:NH1	2.40	0.55
1:A:25:ILE:HG23	1:A:55:LEU:HD13	1.88	0.55
1:A:186:ALA:O	1:A:189:LEU:HB2	2.07	0.55
1:B:181:GLN:HG2	1:B:182:GLN:N	2.22	0.55
1:B:131:LEU:O	1:B:135:GLU:HB2	2.07	0.55
1:A:18:GLN:HA	1:A:50:MET:HE1	1.89	0.54
1:B:42:ALA:HA	1:B:63:ALA:O	2.07	0.54
1:A:133:GLU:HG2	1:A:142:VAL:HG21	1.88	0.54
1:A:94:GLY:O	1:A:99:ARG:HD2	2.08	0.54
1:B:105:THR:H	1:B:108:ILE:CG1	2.22	0.53
1:A:45:PHE:HB3	1:B:46:VAL:HG23	1.89	0.53
1:A:46:VAL:HG23	1:B:78:VAL:HG21	1.90	0.52
1:A:249:THR:O	1:A:250:GLN:HB2	2.08	0.52
1:A:127:ILE:HB	1:A:146:GLN:OE1	2.10	0.52
1:B:24:LEU:HD11	1:B:238:LEU:HA	1.92	0.52
1:B:53:GLU:HG2	1:B:54:ARG:HG2	1.91	0.52
1:A:14:CYS:O	1:B:71:SER:HB3	2.10	0.51
1:B:10:ALA:HB1	1:B:237:SER:HB2	1.92	0.51
1:B:191:ARG:HG3	1:B:203:ARG:HG3	1.92	0.50
1:A:35:HIS:CD2	1:A:35:HIS:H	2.29	0.50
1:A:168:PRO:O	1:A:172:ILE:HG12	2.12	0.49
1:B:40:VAL:HG22	1:B:61:VAL:CG2	2.42	0.49
1:A:191:ARG:HE	1:A:203:ARG:HG2	1.77	0.49
1:B:221:THR:O	1:B:224:GLN:HB2	2.13	0.49
1:A:3:LYS:HE2	1:A:226:ARG:O	2.13	0.49
1:A:242:PHE:HA	1:A:245:ILE:HD12	1.95	0.49
1:A:69:ALA:HA	1:A:80:LEU:HD22	1.95	0.48
1:B:57:HIS:HD2	1:B:59:LYS:N	2.00	0.48
1:A:65:GLN:O	1:A:66:ASN:HB2	2.13	0.47
1:A:107:GLU:HG2	1:A:108:ILE:HG13	1.96	0.47
1:A:99:ARG:HG2	1:A:99:ARG:NH2	2.28	0.47
1:A:10:ALA:CB	1:A:237:SER:HB2	2.45	0.47
1:B:24:LEU:HD22	1:B:28:PHE:CZ	2.50	0.47
1:B:129:GLU:HG2	1:B:139:THR:HA	1.97	0.46
1:A:24:LEU:HA	1:A:24:LEU:HD12	1.75	0.46
1:B:162:VAL:CG1	1:B:163:VAL:N	2.79	0.46
1:A:80:LEU:HB2	1:A:81:PRO:HD3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:LEU:C	1:A:27:LEU:HD23	2.37	0.45
1:B:57:HIS:NE2	1:B:59:LYS:HB2	2.32	0.45
1:B:39:CYS:O	1:B:60:PHE:HA	2.17	0.44
1:B:52:LYS:HE3	1:B:86:PHE:O	2.17	0.44
1:A:191:ARG:CD	1:A:203:ARG:HD2	2.41	0.44
1:B:35:HIS:O	1:B:59:LYS:HE2	2.17	0.44
1:B:57:HIS:CD2	1:B:59:LYS:HB2	2.52	0.44
1:A:80:LEU:N	1:A:81:PRO:CD	2.80	0.44
1:A:44:THR:HG23	2:A:605:HOH:O	2.19	0.43
1:B:195:SER:OG	1:B:203:ARG:HD2	2.19	0.43
1:A:197:LYS:O	1:A:198:ILE:HD13	2.18	0.43
1:A:191:ARG:NH2	1:A:206:LEU:O	2.51	0.43
1:A:65:GLN:HB3	1:B:75:THR:HG23	2.01	0.43
1:B:90:TRP:CZ2	1:B:122:MET:HG2	2.53	0.43
1:A:12:TRP:O	1:A:15:ASN:HB2	2.17	0.43
1:B:18:GLN:HA	1:B:50:MET:HE1	2.00	0.43
1:B:105:THR:H	1:B:108:ILE:HG12	1.84	0.43
1:A:250:GLN:NE2	1:A:250:GLN:CA	2.79	0.42
1:A:40:VAL:HA	1:A:61:VAL:O	2.20	0.42
1:B:144:LEU:HD21	1:B:189:LEU:HD22	2.00	0.42
1:B:138:ARG:O	1:B:142:VAL:HG23	2.20	0.42
1:A:79:SER:HB2	1:A:81:PRO:HD2	2.02	0.42
1:B:144:LEU:HD11	1:B:189:LEU:HD21	2.00	0.42
1:A:185:GLU:O	1:A:189:LEU:HD23	2.19	0.41
1:A:206:LEU:HD22	1:A:207:ARG:N	2.35	0.41
1:B:67:ALA:HB2	1:B:83:LEU:HD11	2.03	0.41
1:A:130:THR:HG22	1:A:133:GLU:CD	2.41	0.41
1:A:203:ARG:HG2	1:A:203:ARG:O	2.16	0.41
1:B:5:GLN:NE2	1:B:207:ARG:NH1	2.61	0.41
1:B:19:GLN:O	1:B:22:SER:HB2	2.20	0.41
1:B:154:LEU:HD12	1:B:154:LEU:HA	1.63	0.41
1:B:191:ARG:HG3	1:B:203:ARG:CG	2.51	0.40
1:B:193:TRP:O	1:B:197:LYS:HB2	2.21	0.40
1:A:77:GLU:HG3	1:B:102:TYR:OH	2.22	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:GLU:OE2	1:A:220:ARG:NH1[2_564]	1.72	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:LYS:CD	1:A:250:GLN:OE1[2_564]	1.75	0.45
1:A:205:GLU:OE1	1:A:220:ARG:NH1[2_564]	1.77	0.43
1:A:205:GLU:CD	1:A:220:ARG:NH1[2_564]	1.93	0.27
1:A:161:LYS:CE	1:A:250:GLN:OE1[2_564]	2.04	0.16

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	247/250 (99%)	229 (93%)	17 (7%)	1 (0%)	34	66
1	B	247/250 (99%)	230 (93%)	16 (6%)	1 (0%)	34	66
All	All	494/500 (99%)	459 (93%)	33 (7%)	2 (0%)	34	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	19	GLN
1	A	104	GLU

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	197/198 (100%)	174 (88%)	23 (12%)	5	16
1	B	197/198 (100%)	177 (90%)	20 (10%)	7	22

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	394/396 (100%)	351 (89%)	43 (11%)	6 19

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	SER
1	A	53	GLU
1	A	54	ARG
1	A	55	LEU
1	A	74	PHE
1	A	80	LEU
1	A	99	ARG
1	A	101	TYR
1	A	107	GLU
1	A	131	LEU
1	A	134	ARG
1	A	136	SER
1	A	143	VAL
1	A	152	LYS
1	A	155	LYS
1	A	162	VAL
1	A	177	VAL
1	A	191	ARG
1	A	197	LYS
1	A	202	VAL
1	A	206	LEU
1	A	207	ARG
1	A	250	GLN
1	B	19	GLN
1	B	53	GLU
1	B	89	ASN
1	B	101	TYR
1	B	129	GLU
1	B	131	LEU
1	B	136	SER
1	B	147	ILE
1	B	153	LYS
1	B	154	LEU
1	B	176	LYS
1	B	181	GLN
1	B	189	LEU
1	B	191	ARG

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Mol	Chain	Res	Type
1	B	192	SER
1	B	195	SER
1	B	203	ARG
1	B	205	GLU
1	B	239	LYS
1	B	247	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	HIS
1	A	38	GLN
1	A	57	HIS
1	A	250	GLN
1	B	5	GLN
1	B	35	HIS
1	B	38	GLN
1	B	57	HIS
1	B	66	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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 ⓘ

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.