



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:14 PM UTC

PDB ID : 3TIM / pdb_00003tim
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 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

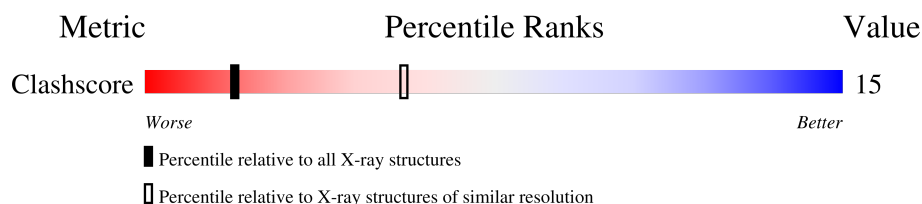
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	190562	4276 (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 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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	250	 63% 33% .
1	B	250	 63% 32% 5%

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		

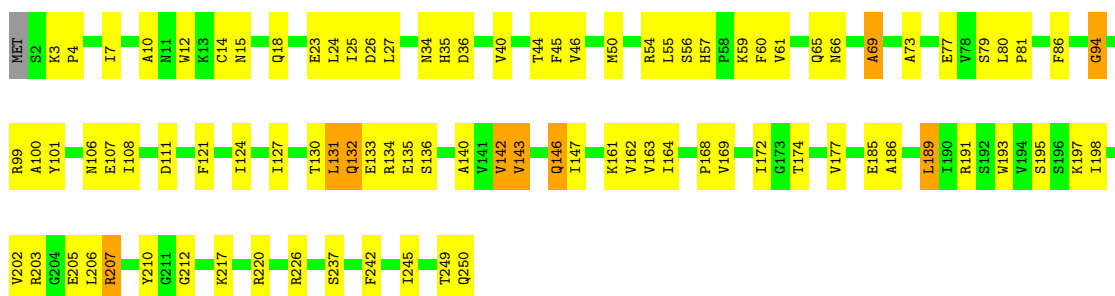
3 Residue-property plots

These plots are drawn for all protein, RNA, 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. 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.

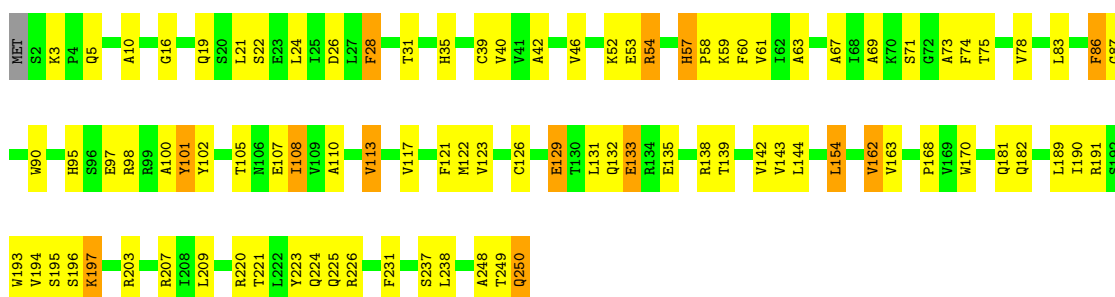
• Molecule 1: TRIOSEPHOSPHATE ISOMERASE

Chain A: 



• Molecule 1: TRIOSEPHOSPHATE ISOMERASE

Chain B: 



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	1.18	3/1923 (0.2%)	1.85	36/2606 (1.4%)
1	B	1.14	4/1923 (0.2%)	1.83	47/2606 (1.8%)
All	All	1.16	7/3846 (0.2%)	1.84	83/5212 (1.6%)

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 (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	168	PRO	CA-C	-6.06	1.46	1.53
1	B	86	PHE	C-O	5.88	1.31	1.24
1	A	163	VAL	CA-C	-5.76	1.45	1.52
1	A	193	TRP	C-O	-5.49	1.17	1.24
1	B	133	GLU	CD-OE2	5.27	1.35	1.25
1	A	94	GLY	C-N	-5.27	1.26	1.33
1	B	26	ASP	CG-OD2	5.09	1.35	1.25

All (83) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	197	LYS	N-CA-C	13.55	126.13	111.36
1	A	36	ASP	CA-CB-CG	8.38	120.98	112.60
1	A	111	ASP	CA-CB-CG	8.36	120.96	112.60
1	A	189	LEU	CB-CA-C	-8.29	97.73	110.92
1	A	34	ASN	CA-CB-CG	8.26	120.86	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	168	PRO	CB-CA-C	-7.81	100.09	112.27
1	A	177	VAL	CB-CA-C	-7.32	100.73	111.19
1	B	57	HIS	N-CA-C	7.25	118.49	109.57
1	A	143	VAL	N-CA-CB	7.21	123.82	110.77
1	B	95	HIS	CA-CB-CG	-7.21	106.59	113.80
1	B	16	GLY	CA-C-O	7.00	126.77	122.22
1	A	106	ASN	CA-CB-CG	6.92	119.52	112.60
1	B	16	GLY	N-CA-C	6.88	118.48	111.95
1	B	58	PRO	N-CA-CB	6.83	110.06	103.51
1	B	86	PHE	CA-C-N	-6.74	111.97	122.69
1	B	86	PHE	C-N-CA	-6.74	111.97	122.69
1	B	28	PHE	N-CA-CB	6.74	119.78	110.01
1	A	60	PHE	N-CA-C	6.34	119.24	108.90
1	B	69	ALA	N-CA-C	6.27	117.80	110.97
1	B	126	CYS	N-CA-C	6.25	119.13	109.07
1	A	202	VAL	N-CA-CB	6.18	118.49	110.57
1	B	196	SER	CA-C-N	6.18	129.06	120.29
1	B	196	SER	C-N-CA	6.18	129.06	120.29
1	B	113	VAL	N-CA-CB	6.08	117.67	110.55
1	B	249	THR	N-CA-CB	5.99	120.96	110.42
1	B	54	ARG	N-CA-C	5.98	120.76	112.45
1	B	162	VAL	N-CA-C	5.97	117.95	108.87
1	B	57	HIS	CB-CA-C	5.97	118.64	109.50
1	B	57	HIS	CA-C-O	5.91	124.67	119.71
1	B	249	THR	N-CA-C	5.91	120.22	113.19
1	B	74	PHE	CA-CB-CG	-5.89	107.91	113.80
1	B	220	ARG	CA-C-N	5.85	128.12	120.28
1	B	220	ARG	C-N-CA	5.85	128.12	120.28
1	B	129	GLU	N-CA-CB	5.85	119.94	110.41
1	A	142	VAL	N-CA-C	5.81	115.94	110.30
1	B	100	ALA	CA-C-N	-5.64	113.62	122.65
1	B	100	ALA	C-N-CA	-5.64	113.62	122.65
1	B	249	THR	CA-C-N	5.63	131.83	121.70
1	B	249	THR	C-N-CA	5.63	131.83	121.70
1	A	207	ARG	NE-CZ-NH1	-5.62	115.88	121.50
1	B	154	LEU	N-CA-C	5.62	119.21	110.17
1	A	132	GLN	CB-CA-C	5.61	119.59	110.96
1	B	250	GLN	CB-CA-C	-5.60	99.46	110.10
1	B	101	TYR	CA-CB-CG	5.60	123.98	113.90
1	A	217	LYS	N-CA-C	5.59	118.22	111.40
1	A	162	VAL	N-CA-C	5.58	116.51	108.48
1	B	108	ILE	CB-CA-C	-5.56	104.86	111.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	121	PHE	N-CA-C	5.55	118.33	110.50
1	A	69	ALA	N-CA-C	5.54	119.76	112.89
1	A	106	ASN	N-CA-CB	5.52	118.16	109.94
1	A	210	TYR	O-C-N	5.50	129.50	123.22
1	B	73	ALA	N-CA-C	5.46	118.07	109.39
1	A	56	SER	CA-C-N	5.45	130.78	121.35
1	A	56	SER	C-N-CA	5.45	130.78	121.35
1	A	100	ALA	N-CA-C	5.42	116.99	111.14
1	B	21	LEU	CA-C-N	5.41	128.30	120.79
1	B	21	LEU	C-N-CA	5.41	128.30	120.79
1	B	101	TYR	N-CA-C	5.40	119.57	112.92
1	A	169	VAL	O-C-N	5.39	127.17	121.89
1	A	7	ILE	CB-CA-C	-5.38	102.60	110.62
1	A	163	VAL	CA-CB-CG2	-5.31	101.37	110.40
1	B	110	ALA	CA-C-N	5.29	127.32	120.44
1	B	110	ALA	C-N-CA	5.29	127.32	120.44
1	A	174	THR	O-C-N	5.28	127.62	122.19
1	B	132	GLN	CB-CG-CD	5.27	121.57	112.60
1	B	31	THR	CA-C-N	-5.26	113.64	122.17
1	B	31	THR	C-N-CA	-5.26	113.64	122.17
1	A	146	GLN	OE1-CD-NE2	5.22	127.82	122.60
1	B	209	LEU	CA-C-N	-5.22	115.75	123.05
1	B	209	LEU	C-N-CA	-5.22	115.75	123.05
1	B	123	VAL	N-CA-C	5.18	116.43	108.71
1	A	134	ARG	NE-CZ-NH2	-5.17	114.54	119.20
1	A	26	ASP	N-CA-CB	5.16	117.80	110.16
1	A	140	ALA	N-CA-CB	5.16	117.70	110.12
1	B	121	PHE	N-CA-CB	5.15	118.70	110.46
1	B	231	PHE	N-CA-C	5.15	117.63	109.50
1	A	131	LEU	CB-CA-C	-5.15	102.25	110.79
1	A	23	GLU	CA-C-N	5.12	127.56	120.29
1	A	23	GLU	C-N-CA	5.12	127.56	120.29
1	A	86	PHE	N-CA-C	5.10	117.58	111.71
1	A	212	GLY	CA-C-N	5.07	129.86	122.41
1	A	212	GLY	C-N-CA	5.07	129.86	122.41
1	A	4	PRO	N-CA-CB	5.00	109.01	102.86

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 [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	1889	0	1925	60	5
1	B	1889	0	1925	62	0
2	A	23	0	0	1	0
2	B	7	0	0	0	0
All	All	3808	0	3850	115	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 (115) 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	0.78
1:A:143:VAL:O	1:A:147:ILE:HG22	1.87	0.74
1:B:5:GLN:HE21	1:B:207:ARG:HH12	1.36	0.72
1:B:52:LYS:HZ1	1:B:87:GLY:HA3	1.53	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:A:73:ALA:HB1	1:B:97:GLU:OE2	1.94	0.68
1:B:105:THR:OG1	1:B:108:ILE:HG12	1.93	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:O	1:B:226:ARG:HG3	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:HD23	1:A:27:LEU:O	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:79:SER:CB	1:A:81:PRO:HD2	2.35	0.56
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:B:133:GLU:HG2	1:B:138:ARG:NH1	2.21	0.56
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: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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:H	1:A:35:HIS:CD2	2.29	0.50
1:A:27:LEU:HD23	1:A:27:LEU:C	2.37	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:170:TRP:H	1:B:170:TRP:CD1	2.29	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:A:250:GLN:NE2	1:A:250:GLN:HA	2.12	0.48
1:B:57:HIS:HD2	1:B:59:LYS:N	2.00	0.48
1:A:18:GLN:HA	1:A:50:MET:HE1	1.95	0.47
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:A:130:THR:HG22	1:A:133:GLU:CD	2.41	0.46
1:B:129:GLU:HG2	1:B:139:THR:HA	1.97	0.46
1:A:24:LEU:HD12	1:A:24:LEU:HA	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
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:191:ARG:NH2	1:A:206:LEU:O	2.51	0.43
1:A:197:LYS:O	1:A:198:ILE:HD13	2.18	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:107:GLU:H	1:B:107:GLU:CD	2.27	0.42
1:B:144:LEU:HD11	1:B:189:LEU:HD21	2.00	0.42
1:B:3:LYS:NZ	1:B:225:GLN:O	2.51	0.41
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: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:HA	1:B:154:LEU:HD12	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:18:GLN:HG2	1:A:50:MET:HE2	2.03	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
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](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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

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.