



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 09:35 AM EDT

PDB ID : 3VNI / pdb_00003vni
Title : Crystal structures of D-Psicose 3-epimerase from *Clostridium cellulolyticum* H10 and its complex with ketohexose sugars
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Deposited on : 2012-01-16
Resolution : 1.98 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

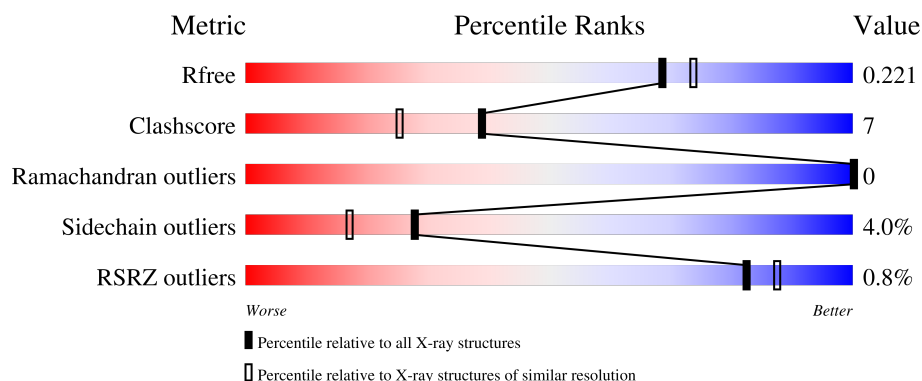
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.98 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	1506 (1.98-1.98)
Clashscore	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)
RSRZ outliers	180081	1506 (1.98-1.98)

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	294	
1	B	294	
1	C	294	
1	D	294	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10344 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 Xylose isomerase domain protein TIM barrel.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2293	1459	388	435	11			
1	B	289	Total	C	N	O	S	0	0	0
			2292	1459	388	435	10			
1	C	288	Total	C	N	O	S	0	0	0
			2287	1456	387	434	10			
1	D	288	Total	C	N	O	S	0	0	0
			2287	1456	387	434	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ALA	-	expression tag	UNP B8I944
B	0	ALA	-	expression tag	UNP B8I944
C	0	ALA	-	expression tag	UNP B8I944
D	0	ALA	-	expression tag	UNP B8I944

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mn	0	0
			1	1		
2	B	1	Total	Mn	0	0
			1	1		
2	C	1	Total	Mn	0	0
			1	1		
2	D	1	Total	Mn	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	339	Total 339	O 339	0	0
3	B	282	Total 282	O 282	0	0
3	C	295	Total 295	O 295	0	0
3	D	265	Total 265	O 265	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.76Å 115.44Å 91.63Å 90.00° 105.45° 90.00°	Depositor
Resolution (Å)	25.00 – 1.98 25.00 – 1.98	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-1.98) 93.8 (25.00-1.98)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.00 (at 1.98Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.175 , 0.220 0.176 , 0.221	Depositor DCC
R_{free} test set	5270 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	22.6	Xtriage
Anisotropy	0.175	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 51.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10344	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 3.84% 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: MN

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	3/2346 (0.1%)	1.07	4/3177 (0.1%)
1	B	0.87	1/2345 (0.0%)	1.06	6/3176 (0.2%)
1	C	0.85	1/2340 (0.0%)	1.06	11/3169 (0.3%)
1	D	0.83	1/2340 (0.0%)	1.06	12/3169 (0.4%)
All	All	0.89	6/9371 (0.1%)	1.06	33/12691 (0.3%)

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	1
1	B	0	1
1	C	0	2
1	D	0	1
All	All	0	5

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	249	MET	SD-CE	-7.79	1.60	1.79
1	A	181	MET	SD-CE	-5.71	1.65	1.79
1	D	224	PRO	CA-C	5.17	1.57	1.52
1	A	64	VAL	C-O	5.11	1.29	1.24
1	A	243	MET	SD-CE	-5.09	1.66	1.79
1	C	249	MET	SD-CE	-5.07	1.66	1.79

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	190	GLU	N-CA-C	8.68	124.00	113.23
1	B	176	ASN	N-CA-C	8.62	125.50	111.37
1	B	190	GLU	N-CA-C	8.13	123.31	113.23
1	C	176	ASN	N-CA-C	7.99	124.55	111.37
1	B	76	SER	N-CA-C	7.54	118.27	109.60
1	D	190	GLU	N-CA-C	7.49	122.52	113.23
1	D	239	GLY	N-CA-C	-7.49	102.48	112.81
1	D	76	SER	N-CA-C	7.45	118.17	109.60
1	D	176	ASN	N-CA-C	7.20	123.25	111.37
1	C	250	GLY	N-CA-C	7.07	121.65	112.18
1	B	250	GLY	N-CA-C	7.06	121.64	112.18
1	D	250	GLY	N-CA-C	6.99	121.54	112.18
1	C	190	GLU	N-CA-C	6.64	121.61	112.90
1	A	286	VAL	N-CA-C	6.57	118.16	111.00
1	C	51	LEU	N-CA-C	-6.55	104.06	111.14
1	D	44	SER	N-CA-C	-6.46	102.19	110.53
1	D	286	VAL	N-CA-C	6.35	117.92	111.00
1	C	286	VAL	N-CA-C	6.11	118.72	111.09
1	C	76	SER	N-CA-C	6.10	117.39	109.64
1	B	286	VAL	N-CA-C	5.95	117.48	111.00
1	C	191	GLU	N-CA-C	5.91	119.53	110.20
1	C	38	SER	N-CA-C	5.59	121.28	113.57
1	B	51	LEU	N-CA-C	-5.51	105.35	111.36
1	D	32	ILE	N-CA-C	5.50	116.65	108.46
1	D	216	LYS	N-CA-C	-5.49	103.45	110.53
1	A	191	GLU	N-CA-C	5.34	118.64	110.20
1	C	248	ARG	N-CA-C	5.19	117.86	109.40
1	D	49	ASN	N-CA-C	5.19	116.94	111.28
1	A	250	GLY	N-CA-C	5.17	119.11	112.18
1	C	38	SER	CA-C-N	-5.17	114.46	120.04
1	C	38	SER	C-N-CA	-5.17	114.46	120.04
1	D	32	ILE	CB-CA-C	-5.05	102.72	110.50
1	D	159	LEU	N-CA-C	5.05	117.57	111.40

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	7	TYR	Sidechain
1	B	7	TYR	Sidechain
1	C	21	TYR	Sidechain
1	C	7	TYR	Sidechain
1	D	7	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	2293	0	2228	32	0
1	B	2292	0	2228	39	0
1	C	2287	0	2223	26	0
1	D	2287	0	2223	28	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	339	0	0	6	1
3	B	282	0	0	1	0
3	C	295	0	0	0	0
3	D	265	0	0	3	1
All	All	10344	0	8902	122	1

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 7.

All (122) 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:150:GLU:HB2	1:B:181:MET:HE2	1.50	0.94
1:C:118:LYS:HA	1:C:118:LYS:HE2	1.57	0.86
1:C:150:GLU:HB2	1:C:181:MET:HE2	1.56	0.85
1:A:75:SER:O	1:A:128:ARG:NH1	2.13	0.81
1:D:123:LYS:HE3	1:D:123:LYS:HA	1.62	0.81
1:D:14:TRP:HE1	1:D:256:ASN:HD22	1.29	0.81
1:D:23:GLU:O	1:D:27:LYS:HG2	1.81	0.80
1:A:98:LYS:HD3	3:A:7402:HOH:O	1.84	0.77
1:D:32:ILE:HD13	1:D:61:THR:HB	1.66	0.77
1:C:157:ASN:HD22	1:C:158:TYR:H	1.33	0.77
1:B:270:LYS:HZ3	1:B:274:ARG:HE	1.34	0.76
1:A:14:TRP:HE1	1:A:256:ASN:HD22	1.35	0.74
1:D:4:GLY:HA3	1:D:32:ILE:HB	1.71	0.72
1:C:25:VAL:HG13	1:C:30:PHE:HB2	1.71	0.72
1:B:274:ARG:CG	1:B:274:ARG:HH11	2.03	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:14:TRP:HE1	1:B:256:ASN:HD22	1.39	0.69
1:B:274:ARG:HH11	1:B:274:ARG:CB	2.08	0.67
1:B:270:LYS:NZ	1:B:274:ARG:HE	1.92	0.66
1:A:1:MET:HE2	1:A:288:GLU:HG2	1.78	0.66
1:D:123:LYS:O	1:D:127:GLU:HG2	1.94	0.66
1:C:14:TRP:HE1	1:C:256:ASN:HD22	1.44	0.66
1:A:271:MET:HE2	1:A:275:GLU:HG2	1.77	0.65
1:A:14:TRP:HE1	1:A:256:ASN:ND2	1.95	0.65
1:B:274:ARG:NH1	1:B:274:ARG:HG2	2.11	0.64
1:C:157:ASN:ND2	1:C:158:TYR:H	1.96	0.64
1:A:115:ASP:OD2	1:A:117:THR:HB	1.98	0.64
1:A:289:CYS:HB3	3:A:7321:HOH:O	1.98	0.63
1:D:14:TRP:HE1	1:D:256:ASN:ND2	1.94	0.63
1:B:14:TRP:HE1	1:B:256:ASN:ND2	1.97	0.62
1:A:66:HIS:HD2	3:A:7421:HOH:O	1.81	0.61
1:D:181:MET:HE1	1:D:209:HIS:CE1	2.36	0.61
1:C:284:ARG:O	1:C:288:GLU:HG3	2.01	0.60
1:C:14:TRP:HE1	1:C:256:ASN:ND2	1.98	0.60
1:B:274:ARG:HH11	1:B:274:ARG:HG2	1.66	0.59
1:C:157:ASN:HD22	1:C:158:TYR:N	2.02	0.57
1:A:150:GLU:HB2	1:A:181:MET:HE2	1.87	0.57
1:D:115:ASP:OD2	1:D:117:THR:HB	2.05	0.56
1:D:249:MET:HE2	1:D:262:ASP:HB2	1.86	0.56
1:B:25:VAL:HG13	1:B:30:PHE:HB2	1.86	0.56
1:A:150:GLU:HB2	1:A:181:MET:CE	2.36	0.55
1:A:271:MET:HE2	1:A:275:GLU:CG	2.36	0.55
1:D:1:MET:HG3	1:D:288:GLU:OE2	2.05	0.55
1:A:1:MET:N	3:A:7183:HOH:O	2.40	0.55
1:A:46:ILE:O	1:A:50:GLU:HG3	2.06	0.55
1:D:181:MET:CE	1:D:209:HIS:CE1	2.89	0.55
1:A:244:GLU:HG2	1:A:246:PHE:CE2	2.43	0.54
1:B:107:ALA:HB2	1:B:152:LEU:HD11	1.89	0.53
3:A:7431:HOH:O	1:D:258:LYS:HE2	2.07	0.53
1:D:42:PHE:CD1	1:D:42:PHE:N	2.72	0.53
1:B:47:GLN:HA	1:B:50:GLU:HG2	1.91	0.53
1:C:150:GLU:HB2	1:C:181:MET:CE	2.35	0.52
1:D:48:ILE:HD11	1:D:95:ARG:HH11	1.74	0.52
1:C:115:ASP:OD2	1:C:118:LYS:HG2	2.10	0.52
1:D:227:GLU:OE1	3:D:644:HOH:O	2.19	0.51
1:B:150:GLU:CB	1:B:181:MET:HE2	2.33	0.51
1:B:249:MET:HE2	1:B:262:ASP:N	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:172:GLN:HG3	3:D:549:HOH:O	2.12	0.50
1:B:212:GLU:CD	1:B:221:GLY:HA3	2.38	0.49
1:B:9:TYR:CD1	1:B:245:PRO:HG2	2.48	0.48
1:A:238:ASN:C	3:A:7420:HOH:O	2.55	0.48
1:B:94:LYS:HD2	1:B:143:CYS:SG	2.53	0.48
1:D:209:HIS:HA	1:D:242:VAL:O	2.14	0.48
1:B:44:SER:H	1:B:47:GLN:CD	2.22	0.47
1:A:5:ILE:HG13	1:A:243:MET:HG3	1.96	0.47
1:C:37:ALA:HB2	1:C:64:VAL:CG1	2.44	0.47
1:C:107:ALA:HB2	1:C:152:LEU:HD11	1.97	0.47
1:B:44:SER:H	1:B:47:GLN:CG	2.27	0.47
1:C:237:TYR:CZ	1:C:239:GLY:HA3	2.50	0.47
1:D:38:SER:O	1:D:41:PRO:HD2	2.15	0.47
1:D:89:TYR:O	1:D:93:LEU:HB2	2.15	0.47
1:D:74:LEU:HD22	1:D:132:SER:HB3	1.96	0.46
1:B:177:ASN:HD22	1:B:177:ASN:N	2.14	0.46
1:D:40:LEU:HB2	1:D:41:PRO:HD3	1.97	0.46
1:C:157:ASN:ND2	1:C:158:TYR:N	2.62	0.46
1:B:21:TYR:O	1:B:25:VAL:HB	2.16	0.46
1:B:260:TRP:CD1	1:C:158:TYR:HA	2.51	0.45
1:A:124:GLY:O	1:A:127:GLU:HG2	2.16	0.45
1:B:270:LYS:NZ	1:B:274:ARG:NE	2.64	0.45
1:B:288:GLU:O	3:B:479:HOH:O	2.21	0.45
1:C:78:ASP:OD1	1:C:78:ASP:C	2.60	0.45
1:B:42:PHE:CD1	1:B:42:PHE:N	2.84	0.45
1:A:271:MET:CE	1:A:275:GLU:HG2	2.45	0.45
1:A:103:LEU:C	1:A:103:LEU:HD23	2.42	0.45
1:A:184:THR:HA	1:A:187:MET:HE3	1.99	0.45
1:A:271:MET:CE	1:A:274:ARG:NH1	2.80	0.45
1:B:116:TYR:CE1	1:C:260:TRP:CZ2	3.05	0.45
1:C:103:LEU:C	1:C:103:LEU:HD23	2.42	0.45
1:A:271:MET:HE1	1:A:274:ARG:NH1	2.32	0.45
1:B:96:LEU:HG	1:B:101:VAL:HG22	1.98	0.45
1:C:118:LYS:HE2	1:C:118:LYS:CA	2.34	0.45
1:A:249:MET:HE2	1:A:262:ASP:HB2	1.97	0.44
1:B:4:GLY:HA3	1:B:32:ILE:HB	1.99	0.44
1:B:157:ASN:ND2	1:B:158:TYR:H	2.15	0.44
1:C:120:ILE:HD12	1:C:158:TYR:CE2	2.53	0.44
1:A:249:MET:HE2	1:A:262:ASP:N	2.32	0.44
1:A:150:GLU:OE1	1:A:152:LEU:HD21	2.18	0.44
1:A:30:PHE:CE1	1:A:280:LEU:HG	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:ARG:HH11	1:B:274:ARG:HB2	1.83	0.44
1:D:21:TYR:O	1:D:25:VAL:HG22	2.18	0.44
1:B:43:TYR:HA	1:B:47:GLN:OE1	2.18	0.43
1:A:86:LYS:NZ	1:A:135:GLU:OE2	2.38	0.43
1:A:74:LEU:HD22	1:A:132:SER:HB3	2.00	0.43
1:C:177:ASN:N	1:C:177:ASN:HD22	2.17	0.43
1:B:244:GLU:HG2	1:B:246:PHE:CE2	2.54	0.43
1:B:176:ASN:HB3	1:B:177:ASN:HD22	1.83	0.42
1:B:26:ALA:HA	1:B:60:ILE:HD11	2.01	0.42
1:D:111:TYR:CE2	1:D:114:ILE:HA	2.54	0.42
1:D:243:MET:C	1:D:245:PRO:HD3	2.44	0.42
1:C:249:MET:HE2	1:C:262:ASP:N	2.35	0.42
1:B:118:LYS:HZ2	1:B:119:THR:H	1.65	0.42
1:B:173:VAL:HG11	1:B:178:VAL:HG21	2.02	0.42
1:B:96:LEU:HG	1:B:101:VAL:CG2	2.50	0.41
1:C:99:LEU:O	1:C:100:ASP:HB3	2.20	0.41
1:D:30:PHE:CE1	1:D:280:LEU:HG	2.55	0.41
1:B:258:LYS:HG3	1:B:260:TRP:CZ2	2.56	0.41
1:C:7:TYR:CG	1:C:8:ALA:N	2.89	0.41
1:A:111:TYR:CE1	1:A:114:ILE:HG12	2.56	0.41
1:D:157:ASN:HB3	3:D:639:HOH:O	2.21	0.41
1:A:120:ILE:HD12	1:A:158:TYR:CE2	2.55	0.41
1:A:193:SER:HB2	1:C:230:GLU:OE2	2.20	0.41
1:B:42:PHE:H	1:B:42:PHE:HD1	1.67	0.40
1:D:103:LEU:C	1:D:103:LEU:HD23	2.45	0.40

All (1) 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)
3:A:7236:HOH:O	3:D:645:HOH:O[2_545]	2.19	0.01

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	287/294 (98%)	281 (98%)	6 (2%)	0	100	100
1	B	287/294 (98%)	278 (97%)	9 (3%)	0	100	100
1	C	286/294 (97%)	280 (98%)	6 (2%)	0	100	100
1	D	286/294 (97%)	279 (98%)	7 (2%)	0	100	100
All	All	1146/1176 (97%)	1118 (98%)	28 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	241/245 (98%)	232 (96%)	9 (4%)	30	20
1	B	240/245 (98%)	230 (96%)	10 (4%)	26	15
1	C	240/245 (98%)	232 (97%)	8 (3%)	33	23
1	D	240/245 (98%)	229 (95%)	11 (5%)	24	12
All	All	961/980 (98%)	923 (96%)	38 (4%)	28	17

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	27	LYS
1	A	62	LEU
1	A	96	LEU
1	A	116	TYR
1	A	119	THR
1	A	272	LEU
1	A	280	LEU
1	A	289	CYS
1	B	25	VAL

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Mol	Chain	Res	Type
1	B	42	PHE
1	B	47	GLN
1	B	96	LEU
1	B	101	VAL
1	B	116	TYR
1	B	118	LYS
1	B	177	ASN
1	B	205	LEU
1	B	274	ARG
1	C	1	MET
1	C	25	VAL
1	C	62	LEU
1	C	96	LEU
1	C	99	LEU
1	C	118	LYS
1	C	270	LYS
1	C	272	LEU
1	D	2	LYS
1	D	19	LYS
1	D	42	PHE
1	D	62	LEU
1	D	93	LEU
1	D	96	LEU
1	D	99	LEU
1	D	116	TYR
1	D	118	LYS
1	D	123	LYS
1	D	280	LEU

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	66	HIS
1	A	256	ASN
1	A	265	ASN
1	A	277	GLN
1	B	157	ASN
1	B	176	ASN
1	B	177	ASN
1	B	256	ASN
1	C	49	ASN
1	C	66	HIS

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Mol	Chain	Res	Type
1	C	84	ASN
1	C	157	ASN
1	C	177	ASN
1	C	238	ASN
1	C	256	ASN
1	D	177	ASN
1	D	256	ASN
1	D	265	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](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

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	289/294 (98%)	-0.50	1 (0%) 90 93	16, 23, 38, 58	0
1	B	289/294 (98%)	-0.26	5 (1%) 69 78	17, 27, 44, 58	0
1	C	288/294 (97%)	-0.27	0 100 100	17, 28, 42, 62	0
1	D	288/294 (97%)	-0.08	3 (1%) 79 86	17, 30, 47, 61	0
All	All	1154/1176 (98%)	-0.28	9 (0%) 82 87	16, 27, 44, 62	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	42	PHE	5.1
1	B	0	ALA	3.2
1	B	42	PHE	3.0
1	B	118	LYS	2.9
1	A	289	CYS	2.4
1	D	48	ILE	2.0
1	D	43	TYR	2.0
1	B	119	THR	2.0
1	B	265	ASN	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	MN	C	301	1/1	0.96	0.10	44,44,44,44	0
2	MN	B	301	1/1	0.98	0.08	37,37,37,37	0
2	MN	A	7001	1/1	0.99	0.12	41,41,41,41	0
2	MN	D	301	1/1	0.99	0.09	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.