



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 03:54 AM GMT

PDB ID : 1PW4
Title : Crystal Structure of the Glycerol-3-Phosphate Transporter from E.Coli
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Deposited on : 2003-06-30
Resolution : 3.30 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

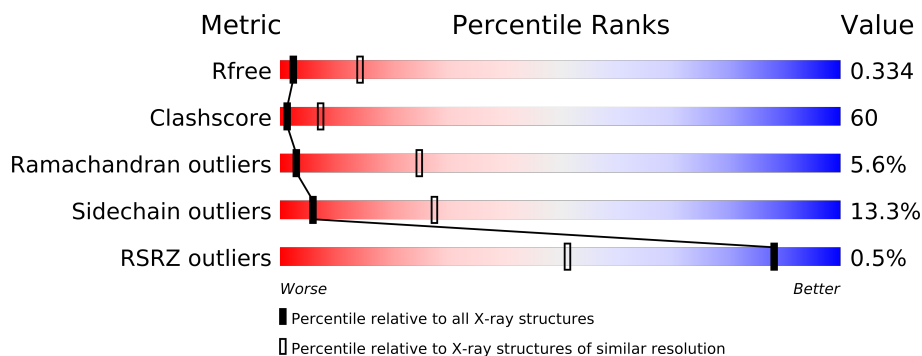
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.30 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}	66092	1341 (3.40-3.20)
Clashscore	79885	1696 (3.40-3.20)
Ramachandran outliers	78287	1664 (3.40-3.20)
Sidechain outliers	78261	1662 (3.40-3.20)
RSRZ outliers	66119	1342 (3.40-3.20)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	451	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3403 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Glycerol-3-phosphate transporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	434	3403	2275	546	554	28	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

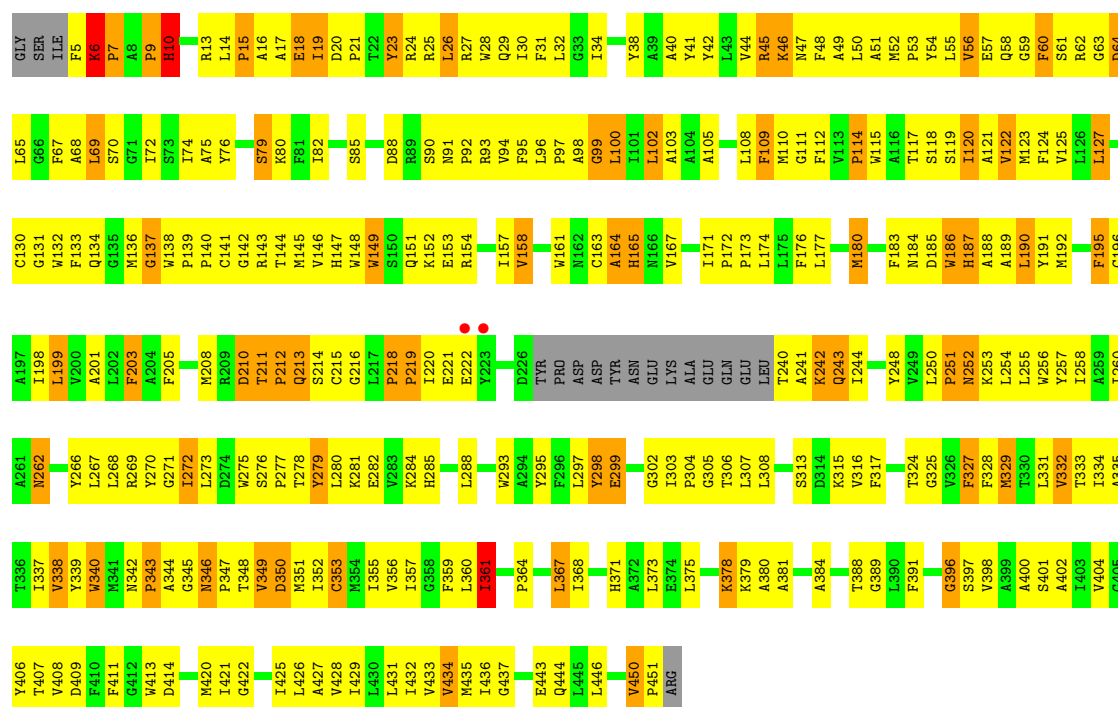
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	GLY	-	CLONING ARTIFACT	UNP P08194
A	449	LEU	-	CLONING ARTIFACT	UNP P08194
A	450	VAL	-	CLONING ARTIFACT	UNP P08194
A	451	PRO	-	CLONING ARTIFACT	UNP P08194
A	452	ARG	-	CLONING ARTIFACT	UNP P08194

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 errors 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: Glycerol-3-phosphate transporter

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	97.64Å 97.64Å 175.15Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 3.30 14.98 – 3.30	Depositor EDS
% Data completeness (in resolution range)	95.4 (15.00-3.30) 95.4 (14.98-3.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.84 (at 3.32Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.296 , 0.325 0.306 , 0.334	Depositor DCC
R_{free} test set	720 reflections (5.06%)	DCC
Wilson B-factor (Å ²)	86.1	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 60.9	EDS
Estimated twinning fraction	0.036 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 14843 reflections	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	3403	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.06% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.58	2/3515 (0.1%)	0.95	19/4781 (0.4%)

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	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	164	ALA	C-N	15.29	1.69	1.34
1	A	165	HIS	C-N	-8.11	1.15	1.34

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	45	ARG	C-N-CA	-13.96	86.80	121.70
1	A	164	ALA	C-N-CA	-9.02	99.16	121.70
1	A	57	GLU	N-CA-C	-8.89	87.00	111.00
1	A	63	GLY	N-CA-C	-8.39	92.13	113.10
1	A	165	HIS	O-C-N	-7.87	110.11	122.70
1	A	122	VAL	N-CA-C	-7.73	90.14	111.00
1	A	6	LYS	C-N-CD	-7.28	104.58	120.60
1	A	450	VAL	N-CA-C	6.53	128.62	111.00
1	A	46	LYS	CB-CA-C	6.50	123.39	110.40
1	A	282	GLU	N-CA-C	-5.86	95.19	111.00
1	A	361	ILE	CB-CA-C	-5.80	100.01	111.60
1	A	45	ARG	CA-C-N	5.62	129.56	117.20
1	A	250	LEU	C-N-CD	5.59	140.15	128.40
1	A	218	PRO	N-CA-C	-5.38	98.10	112.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	213	GLN	N-CA-C	5.37	125.49	111.00
1	A	165	HIS	CA-C-N	5.30	128.86	117.20
1	A	199	LEU	CA-CB-CG	5.23	127.33	115.30
1	A	6	LYS	C-N-CA	5.22	143.93	122.00
1	A	45	ARG	O-C-N	-5.19	114.39	122.70

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	10	HIS	Mainchain
1	A	298	TYR	Sidechain
1	A	99	GLY	Mainchain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3403	0	3421	406	0
All	All	3403	0	3421	406	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 60.

All (406) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:164:ALA:C	1:A:165:HIS:N	1.69	1.46
1:A:373:LEU:HD23	1:A:381:ALA:HB1	1.34	1.10
1:A:350:ASP:HA	1:A:353:CYS:HB3	1.38	1.04
1:A:7:PRO:HG3	1:A:90:SER:HA	1.38	1.00
1:A:281:LYS:HA	1:A:285:HIS:HB3	1.45	0.98
1:A:55:LEU:HD22	1:A:56:VAL:HG13	1.47	0.96
1:A:157:ILE:HG22	1:A:161:TRP:CD1	2.02	0.93
1:A:55:LEU:O	1:A:56:VAL:HG22	1.68	0.93
1:A:13:ARG:HG2	1:A:15:PRO:HD3	1.52	0.90

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:157:ILE:HG22	1:A:161:TRP:NE1	1.87	0.90
1:A:164:ALA:C	1:A:165:HIS:CA	2.39	0.89
1:A:255:LEU:HD12	1:A:258:ILE:HD11	1.55	0.89
1:A:55:LEU:HD22	1:A:56:VAL:CG1	2.02	0.89
1:A:145:MET:SD	1:A:157:ILE:HD13	2.12	0.89
1:A:56:VAL:HG12	1:A:60:PHE:HB3	1.55	0.89
1:A:186:TRP:HD1	1:A:187:HIS:H	1.20	0.88
1:A:275:TRP:HZ3	1:A:400:ALA:O	1.56	0.88
1:A:67:PHE:O	1:A:70:SER:HB3	1.75	0.87
1:A:275:TRP:CZ3	1:A:400:ALA:O	2.28	0.87
1:A:351:MET:O	1:A:355:ILE:HG13	1.74	0.86
1:A:157:ILE:CG2	1:A:161:TRP:NE1	2.39	0.86
1:A:267:LEU:HD13	1:A:400:ALA:HA	1.57	0.86
1:A:62:ARG:HE	1:A:64:ASP:HB2	1.40	0.86
1:A:303:ILE:HB	1:A:304:PRO:HD3	1.58	0.85
1:A:45:ARG:C	1:A:46:LYS:HG2	1.98	0.84
1:A:302:GLY:O	1:A:306:THR:HG22	1.77	0.84
1:A:44:VAL:HA	1:A:47:ASN:HD21	1.39	0.84
1:A:353:CYS:O	1:A:357:ILE:HG12	1.76	0.83
1:A:157:ILE:CG2	1:A:161:TRP:HE1	1.91	0.83
1:A:54:TYR:OH	1:A:288:LEU:HD21	1.79	0.81
1:A:281:LYS:HA	1:A:285:HIS:CB	2.10	0.81
1:A:17:ALA:O	1:A:18:GLU:HB2	1.81	0.80
1:A:154:ARG:O	1:A:158:VAL:HG23	1.82	0.80
1:A:56:VAL:HG11	1:A:65:LEU:HD11	1.64	0.79
1:A:53:PRO:HB3	1:A:278:THR:HA	1.65	0.79
1:A:102:LEU:HD22	1:A:133:PHE:CE1	2.17	0.78
1:A:149:TRP:CD2	1:A:152:LYS:HG3	2.19	0.78
1:A:313:SER:HB2	1:A:324:THR:HG21	1.64	0.78
1:A:198:ILE:HD12	1:A:199:LEU:N	1.99	0.78
1:A:7:PRO:HG3	1:A:90:SER:CA	2.11	0.78
1:A:20:ASP:HB2	1:A:21:PRO:HD3	1.64	0.77
1:A:69:LEU:HD23	1:A:72:ILE:HD13	1.66	0.77
1:A:269:ARG:HG3	1:A:361:ILE:HD12	1.67	0.77
1:A:56:VAL:HG11	1:A:61:SER:HB3	1.67	0.76
1:A:373:LEU:CD2	1:A:381:ALA:HB1	2.14	0.76
1:A:340:TRP:HH2	1:A:414:ASP:OD2	1.67	0.76
1:A:272:ILE:HG13	1:A:420:MET:CE	2.17	0.74
1:A:334:ILE:O	1:A:338:VAL:HG12	1.88	0.74
1:A:41:TYR:HB3	1:A:134:GLN:HE21	1.53	0.74
1:A:96:LEU:HB3	1:A:97:PRO:HD3	1.70	0.74
1:A:328:PHE:O	1:A:332:VAL:HG13	1.87	0.74

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:165:HIS:ND1	1:A:165:HIS:O	2.20	0.73
1:A:339:TYR:CD2	1:A:353:CYS:HB2	2.24	0.73
1:A:163:CYS:O	1:A:167:VAL:HG23	1.88	0.73
1:A:7:PRO:CG	1:A:90:SER:HA	2.15	0.73
1:A:171:ILE:O	1:A:174:LEU:HB2	1.89	0.72
1:A:349:VAL:O	1:A:352:ILE:HG22	1.90	0.72
1:A:6:LYS:HD2	1:A:6:LYS:N	2.04	0.72
1:A:119:SER:HB3	1:A:122:VAL:HB	1.71	0.71
1:A:350:ASP:CA	1:A:353:CYS:HB3	2.19	0.70
1:A:123:MET:O	1:A:127:LEU:HB2	1.91	0.70
1:A:165:HIS:CE1	1:A:299:GLU:OE2	2.44	0.70
1:A:145:MET:SD	1:A:157:ILE:CD1	2.80	0.70
1:A:267:LEU:HD12	1:A:267:LEU:O	1.91	0.70
1:A:157:ILE:HG21	1:A:161:TRP:HE1	1.54	0.69
1:A:284:LYS:HE2	1:A:342:ASN:HB3	1.74	0.69
1:A:13:ARG:HG2	1:A:15:PRO:CD	2.22	0.69
1:A:422:GLY:HA2	1:A:425:ILE:HD12	1.73	0.69
1:A:337:ILE:HG22	1:A:421:ILE:HD13	1.74	0.69
1:A:240:THR:HG22	1:A:241:ALA:H	1.57	0.69
1:A:56:VAL:HG12	1:A:60:PHE:CB	2.22	0.69
1:A:262:ASN:ND2	1:A:368:ILE:HD12	2.07	0.69
1:A:339:TYR:CE1	1:A:353:CYS:HA	2.28	0.68
1:A:346:ASN:HD21	1:A:350:ASP:CG	1.96	0.68
1:A:267:LEU:CD1	1:A:400:ALA:HA	2.24	0.68
1:A:40:ALA:CB	1:A:100:LEU:HD11	2.23	0.68
1:A:220:ILE:O	1:A:222:GLU:N	2.27	0.68
1:A:298:TYR:CZ	1:A:361:ILE:HD11	2.29	0.68
1:A:269:ARG:HB2	1:A:361:ILE:HG21	1.76	0.68
1:A:279:TYR:HD1	1:A:413:TRP:NE1	1.93	0.67
1:A:49:ALA:HA	1:A:52:MET:CE	2.25	0.67
1:A:255:LEU:O	1:A:258:ILE:HG13	1.95	0.66
1:A:337:ILE:O	1:A:340:TRP:HB3	1.95	0.66
1:A:255:LEU:HD21	1:A:384:ALA:HB1	1.76	0.66
1:A:41:TYR:HB3	1:A:134:GLN:NE2	2.11	0.66
1:A:149:TRP:CE2	1:A:152:LYS:HG3	2.31	0.65
1:A:284:LYS:HD3	1:A:346:ASN:ND2	2.12	0.65
1:A:103:ALA:HB2	1:A:133:PHE:HB3	1.79	0.65
1:A:50:LEU:HD11	1:A:173:PRO:HA	1.79	0.65
1:A:72:ILE:HD12	1:A:72:ILE:H	1.62	0.65
1:A:273:LEU:HD22	1:A:298:TYR:CE2	2.33	0.64
1:A:350:ASP:HA	1:A:353:CYS:CB	2.23	0.64
1:A:98:ALA:O	1:A:102:LEU:HB2	1.97	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:258:ILE:HD11	1:A:375:LEU:HD12	1.79	0.64
1:A:210:ASP:HB2	1:A:212:PRO:HD2	1.78	0.64
1:A:40:ALA:HB3	1:A:100:LEU:HD11	1.79	0.64
1:A:138:TRP:O	1:A:141:CYS:HB3	1.98	0.64
1:A:95:PHE:CE2	1:A:136:MET:HB2	2.33	0.63
1:A:272:ILE:HG13	1:A:420:MET:HE3	1.81	0.63
1:A:340:TRP:CE3	1:A:340:TRP:O	2.52	0.63
1:A:364:PRO:O	1:A:368:ILE:HG13	1.97	0.63
1:A:346:ASN:N	1:A:347:PRO:CD	2.61	0.63
1:A:273:LEU:HD22	1:A:298:TYR:HE2	1.64	0.62
1:A:49:ALA:HA	1:A:52:MET:HE3	1.81	0.62
1:A:64:ASP:HA	1:A:406:TYR:CE1	2.35	0.62
1:A:64:ASP:HA	1:A:406:TYR:HE1	1.63	0.62
1:A:52:MET:O	1:A:56:VAL:HG22	1.98	0.62
1:A:44:VAL:HA	1:A:47:ASN:ND2	2.13	0.62
1:A:53:PRO:HB3	1:A:278:THR:CA	2.30	0.62
1:A:60:PHE:CZ	1:A:118:SER:HB3	2.35	0.62
1:A:52:MET:HG2	1:A:65:LEU:HD13	1.81	0.62
1:A:211:THR:N	1:A:212:PRO:CD	2.63	0.62
1:A:191:TYR:CD2	1:A:192:MET:N	2.68	0.62
1:A:195:PHE:O	1:A:198:ILE:HG13	2.00	0.61
1:A:55:LEU:HD23	1:A:56:VAL:H	1.64	0.61
1:A:212:PRO:HG2	1:A:213:GLN:H	1.65	0.61
1:A:103:ALA:HB2	1:A:133:PHE:CB	2.29	0.61
1:A:102:LEU:HD13	1:A:133:PHE:CE2	2.35	0.61
1:A:45:ARG:O	1:A:46:LYS:HG2	2.00	0.61
1:A:103:ALA:HA	1:A:133:PHE:HD1	1.64	0.61
1:A:281:LYS:O	1:A:281:LYS:HG2	2.01	0.61
1:A:408:VAL:HG11	1:A:413:TRP:HE3	1.66	0.61
1:A:298:TYR:OH	1:A:361:ILE:HD11	2.01	0.61
1:A:281:LYS:CA	1:A:285:HIS:HB3	2.27	0.61
1:A:15:PRO:O	1:A:19:ILE:HG12	2.01	0.60
1:A:99:GLY:HA2	1:A:136:MET:SD	2.41	0.60
1:A:408:VAL:HG11	1:A:413:TRP:CE3	2.36	0.60
1:A:220:ILE:C	1:A:222:GLU:H	2.03	0.60
1:A:9:PRO:HG2	1:A:10:HIS:H	1.66	0.60
1:A:298:TYR:OH	1:A:361:ILE:CD1	2.50	0.60
1:A:55:LEU:HD23	1:A:56:VAL:N	2.16	0.60
1:A:94:VAL:HG22	1:A:205:PHE:HE1	1.64	0.60
1:A:275:TRP:HD1	1:A:278:THR:HG21	1.67	0.59
1:A:99:GLY:CA	1:A:136:MET:SD	2.90	0.59
1:A:198:ILE:HD11	1:A:199:LEU:HD23	1.84	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:407:THR:O	1:A:407:THR:HG22	2.03	0.59
1:A:279:TYR:CD1	1:A:413:TRP:NE1	2.71	0.58
1:A:327:PHE:C	1:A:327:PHE:CD1	2.76	0.58
1:A:82:ILE:HD11	1:A:132:TRP:CZ2	2.38	0.58
1:A:211:THR:N	1:A:212:PRO:HD2	2.18	0.58
1:A:210:ASP:C	1:A:212:PRO:CD	2.72	0.58
1:A:396:GLY:O	1:A:400:ALA:HB2	2.04	0.58
1:A:272:ILE:HG13	1:A:420:MET:HE1	1.86	0.58
1:A:10:HIS:ND1	1:A:10:HIS:O	2.36	0.58
1:A:275:TRP:CH2	1:A:401:SER:HA	2.39	0.58
1:A:384:ALA:O	1:A:388:THR:HG23	2.04	0.58
1:A:108:LEU:HD21	1:A:195:PHE:HD1	1.68	0.58
1:A:198:ILE:CD1	1:A:199:LEU:HD23	2.34	0.58
1:A:165:HIS:HD1	1:A:165:HIS:C	2.07	0.57
1:A:56:VAL:CG1	1:A:65:LEU:HD11	2.33	0.57
1:A:339:TYR:CG	1:A:353:CYS:HB2	2.40	0.57
1:A:340:TRP:CH2	1:A:414:ASP:HA	2.39	0.57
1:A:271:GLY:O	1:A:275:TRP:HE3	1.87	0.57
1:A:97:PRO:HG2	1:A:205:PHE:HB2	1.86	0.57
1:A:157:ILE:O	1:A:158:VAL:C	2.42	0.57
1:A:109:PHE:HD2	1:A:109:PHE:O	1.86	0.57
1:A:258:ILE:CD1	1:A:375:LEU:HD12	2.35	0.57
1:A:450:VAL:O	1:A:451:PRO:C	2.42	0.57
1:A:191:TYR:HD2	1:A:192:MET:N	2.03	0.57
1:A:180:MET:HA	1:A:180:MET:CE	2.35	0.57
1:A:45:ARG:O	1:A:46:LYS:CG	2.53	0.56
1:A:171:ILE:N	1:A:172:PRO:HD2	2.20	0.56
1:A:203:PHE:C	1:A:203:PHE:CD2	2.78	0.56
1:A:335:ALA:O	1:A:338:VAL:HG13	2.05	0.56
1:A:42:TYR:HE2	1:A:45:ARG:CZ	2.17	0.56
1:A:56:VAL:HG23	1:A:56:VAL:O	2.06	0.56
1:A:102:LEU:HD22	1:A:133:PHE:CZ	2.39	0.56
1:A:210:ASP:C	1:A:212:PRO:HD3	2.26	0.56
1:A:255:LEU:HD13	1:A:375:LEU:HB2	1.87	0.55
1:A:122:VAL:CG1	1:A:123:MET:N	2.69	0.55
1:A:186:TRP:HD1	1:A:187:HIS:N	1.98	0.55
1:A:428:VAL:O	1:A:432:ILE:HG12	2.07	0.55
1:A:337:ILE:HG13	1:A:338:VAL:N	2.22	0.55
1:A:262:ASN:CG	1:A:368:ILE:HD12	2.27	0.55
1:A:212:PRO:CG	1:A:213:GLN:H	2.20	0.55
1:A:429:ILE:O	1:A:432:ILE:HB	2.08	0.54
1:A:14:LEU:N	1:A:15:PRO:HD3	2.22	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:108:LEU:HD21	1:A:195:PHE:CD1	2.43	0.54
1:A:244:ILE:CD1	1:A:379:LYS:HD3	2.37	0.54
1:A:70:SER:O	1:A:74:ILE:HG13	2.07	0.54
1:A:325:GLY:O	1:A:329:MET:N	2.26	0.54
1:A:72:ILE:HD11	1:A:127:LEU:HG	1.88	0.54
1:A:50:LEU:O	1:A:53:PRO:HD2	2.08	0.54
1:A:218:PRO:O	1:A:220:ILE:N	2.41	0.54
1:A:14:LEU:O	1:A:15:PRO:C	2.46	0.54
1:A:183:PHE:HB3	1:A:191:TYR:HH	1.72	0.54
1:A:346:ASN:N	1:A:347:PRO:HD3	2.23	0.54
1:A:343:PRO:C	1:A:345:GLY:H	2.11	0.54
1:A:211:THR:HG22	1:A:211:THR:O	2.08	0.54
1:A:186:TRP:CD1	1:A:187:HIS:N	2.74	0.54
1:A:436:ILE:CG2	1:A:437:GLY:N	2.71	0.54
1:A:305:GLY:O	1:A:308:LEU:N	2.41	0.54
1:A:27:ARG:NH2	1:A:149:TRP:CE2	2.76	0.53
1:A:119:SER:HB3	1:A:122:VAL:CB	2.39	0.53
1:A:121:ALA:HA	1:A:124:PHE:H	1.73	0.53
1:A:13:ARG:HG3	1:A:216:GLY:HA3	1.90	0.53
1:A:111:GLY:HA2	1:A:190:LEU:HD21	1.90	0.53
1:A:183:PHE:O	1:A:185:ASP:N	2.40	0.53
1:A:275:TRP:CH2	1:A:401:SER:O	2.62	0.53
1:A:183:PHE:HB3	1:A:191:TYR:OH	2.08	0.53
1:A:93:ARG:HA	1:A:208:MET:HE2	1.89	0.53
1:A:58:GLN:O	1:A:60:PHE:N	2.42	0.53
1:A:41:TYR:OH	1:A:137:GLY:HA2	2.09	0.53
1:A:339:TYR:CE2	1:A:353:CYS:HB2	2.44	0.53
1:A:62:ARG:HB2	1:A:64:ASP:HB3	1.90	0.53
1:A:76:TYR:O	1:A:79:SER:HB3	2.09	0.53
1:A:240:THR:HG22	1:A:241:ALA:N	2.23	0.53
1:A:55:LEU:O	1:A:56:VAL:CG2	2.51	0.52
1:A:56:VAL:HB	1:A:61:SER:H	1.74	0.52
1:A:275:TRP:HH2	1:A:401:SER:O	1.92	0.52
1:A:40:ALA:HB1	1:A:100:LEU:HD11	1.91	0.52
1:A:275:TRP:HD1	1:A:278:THR:CG2	2.23	0.52
1:A:275:TRP:CD1	1:A:278:THR:HG21	2.45	0.52
1:A:119:SER:HB3	1:A:122:VAL:CG1	2.40	0.52
1:A:50:LEU:HD11	1:A:173:PRO:CA	2.39	0.52
1:A:327:PHE:HE1	1:A:331:LEU:HD12	1.75	0.52
1:A:45:ARG:CB	1:A:46:LYS:HG2	2.40	0.52
1:A:345:GLY:C	1:A:347:PRO:HD2	2.30	0.52
1:A:212:PRO:O	1:A:216:GLY:N	2.42	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:380:ALA:O	1:A:384:ALA:N	2.42	0.52
1:A:55:LEU:HD22	1:A:56:VAL:HG12	1.88	0.52
1:A:127:LEU:O	1:A:130:CYS:HB3	2.10	0.52
1:A:82:ILE:HD11	1:A:132:TRP:HZ2	1.75	0.52
1:A:198:ILE:O	1:A:201:ALA:HB3	2.10	0.51
1:A:348:THR:HG22	1:A:348:THR:O	2.10	0.51
1:A:56:VAL:HG21	1:A:61:SER:HB3	1.91	0.51
1:A:41:TYR:CE2	1:A:137:GLY:HA3	2.45	0.51
1:A:280:LEU:O	1:A:281:LYS:HB3	2.11	0.51
1:A:23:TYR:HE2	1:A:211:THR:CG2	2.24	0.51
1:A:324:THR:HB	1:A:367:LEU:HD21	1.93	0.51
1:A:99:GLY:HA3	1:A:136:MET:SD	2.51	0.51
1:A:157:ILE:HG22	1:A:161:TRP:HE1	1.59	0.51
1:A:105:ALA:O	1:A:108:LEU:HB2	2.11	0.51
1:A:10:HIS:CG	1:A:10:HIS:O	2.64	0.51
1:A:327:PHE:CE1	1:A:331:LEU:HD12	2.45	0.51
1:A:352:ILE:O	1:A:356:VAL:HG23	2.11	0.51
1:A:52:MET:O	1:A:55:LEU:O	2.28	0.51
1:A:327:PHE:C	1:A:327:PHE:HD1	2.13	0.51
1:A:315:LYS:HD2	1:A:315:LYS:O	2.11	0.51
1:A:335:ALA:O	1:A:339:TYR:CD1	2.64	0.50
1:A:18:GLU:C	1:A:21:PRO:HD2	2.32	0.50
1:A:48:PHE:O	1:A:51:ALA:HB3	2.12	0.50
1:A:55:LEU:CD2	1:A:56:VAL:HG13	2.29	0.50
1:A:213:GLN:OE1	1:A:214:SER:HB2	2.11	0.50
1:A:251:PRO:O	1:A:256:TRP:CD1	2.65	0.50
1:A:94:VAL:O	1:A:97:PRO:HD2	2.11	0.50
1:A:325:GLY:O	1:A:329:MET:HB2	2.11	0.50
1:A:13:ARG:HG2	1:A:15:PRO:CG	2.41	0.50
1:A:429:ILE:O	1:A:433:VAL:HG23	2.11	0.50
1:A:257:TYR:HA	1:A:260:ILE:HG22	1.93	0.50
1:A:218:PRO:O	1:A:219:PRO:C	2.49	0.50
1:A:6:LYS:HD2	1:A:6:LYS:H	1.77	0.50
1:A:242:LYS:O	1:A:243:GLN:C	2.50	0.50
1:A:317:PHE:CE1	1:A:324:THR:HA	2.46	0.50
1:A:157:ILE:HG21	1:A:161:TRP:NE1	2.21	0.49
1:A:212:PRO:HG2	1:A:213:GLN:N	2.27	0.49
1:A:91:ASN:OD1	1:A:210:ASP:OD2	2.30	0.49
1:A:62:ARG:HE	1:A:64:ASP:CB	2.16	0.49
1:A:38:TYR:CE1	1:A:138:TRP:HB2	2.48	0.49
1:A:252:ASN:O	1:A:254:LEU:N	2.46	0.49
1:A:275:TRP:HA	1:A:278:THR:CG2	2.42	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:277:PRO:HG2	1:A:278:THR:H	1.78	0.49
1:A:339:TYR:HE2	1:A:349:VAL:HG12	1.78	0.49
1:A:431:LEU:O	1:A:434:VAL:HG23	2.12	0.49
1:A:284:LYS:HD3	1:A:346:ASN:CB	2.43	0.49
1:A:324:THR:HB	1:A:367:LEU:CD2	2.42	0.49
1:A:13:ARG:C	1:A:15:PRO:HD3	2.33	0.48
1:A:144:THR:O	1:A:147:HIS:N	2.41	0.48
1:A:120:ILE:O	1:A:121:ALA:HB3	2.13	0.48
1:A:51:ALA:O	1:A:55:LEU:N	2.41	0.48
1:A:443:GLU:O	1:A:446:LEU:HB2	2.13	0.48
1:A:211:THR:HA	1:A:213:GLN:HE21	1.78	0.48
1:A:27:ARG:NH2	1:A:149:TRP:NE1	2.62	0.48
1:A:279:TYR:C	1:A:279:TYR:CD2	2.86	0.48
1:A:210:ASP:CB	1:A:212:PRO:HD2	2.44	0.48
1:A:7:PRO:HG3	1:A:90:SER:CB	2.42	0.47
1:A:20:ASP:HB3	1:A:24:ARG:HH21	1.78	0.47
1:A:284:LYS:HD3	1:A:346:ASN:HD22	1.79	0.47
1:A:267:LEU:HD13	1:A:400:ALA:CA	2.37	0.47
1:A:285:HIS:ND1	1:A:285:HIS:N	2.60	0.47
1:A:64:ASP:O	1:A:67:PHE:N	2.46	0.47
1:A:49:ALA:HA	1:A:52:MET:HE2	1.95	0.47
1:A:91:ASN:HA	1:A:92:PRO:HD3	1.77	0.47
1:A:40:ALA:HB3	1:A:100:LEU:CD1	2.45	0.47
1:A:14:LEU:N	1:A:15:PRO:CD	2.77	0.47
1:A:268:LEU:O	1:A:268:LEU:HD13	2.14	0.47
1:A:20:ASP:HB2	1:A:21:PRO:CD	2.40	0.47
1:A:28:TRP:O	1:A:32:LEU:HD13	2.15	0.47
1:A:165:HIS:ND1	1:A:165:HIS:C	2.67	0.46
1:A:45:ARG:C	1:A:46:LYS:CG	2.68	0.46
1:A:112:PHE:O	1:A:114:PRO:HD3	2.15	0.46
1:A:198:ILE:C	1:A:198:ILE:HD12	2.34	0.46
1:A:55:LEU:CD2	1:A:56:VAL:CG1	2.86	0.46
1:A:23:TYR:HE2	1:A:211:THR:HG21	1.79	0.46
1:A:345:GLY:C	1:A:347:PRO:CD	2.84	0.46
1:A:421:ILE:O	1:A:425:ILE:HG13	2.15	0.46
1:A:18:GLU:OE1	1:A:21:PRO:HG2	2.15	0.46
1:A:41:TYR:CZ	1:A:137:GLY:CA	2.99	0.46
1:A:293:TRP:O	1:A:297:LEU:HG	2.14	0.46
1:A:110:MET:O	1:A:117:THR:HB	2.16	0.46
1:A:56:VAL:CG1	1:A:61:SER:H	2.28	0.46
1:A:210:ASP:N	1:A:210:ASP:OD1	2.48	0.46
1:A:65:LEU:O	1:A:68:ALA:HB3	2.15	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:13:ARG:CG	1:A:15:PRO:HD3	2.36	0.46
1:A:133:PHE:O	1:A:136:MET:HG2	2.16	0.46
1:A:268:LEU:O	1:A:272:ILE:HB	2.16	0.46
1:A:102:LEU:HD13	1:A:133:PHE:CZ	2.50	0.46
1:A:136:MET:O	1:A:139:PRO:HD2	2.16	0.46
1:A:404:VAL:O	1:A:408:VAL:HG23	2.15	0.46
1:A:444:GLN:C	1:A:446:LEU:N	2.70	0.46
1:A:346:ASN:ND2	1:A:350:ASP:HB2	2.30	0.46
1:A:34:ILE:HD13	1:A:144:THR:HG21	1.98	0.46
1:A:244:ILE:HD11	1:A:379:LYS:HD3	1.97	0.45
1:A:328:PHE:CD2	1:A:360:LEU:HD22	2.51	0.45
1:A:340:TRP:CH2	1:A:414:ASP:OD2	2.58	0.45
1:A:120:ILE:H	1:A:123:MET:CE	2.29	0.45
1:A:190:LEU:HD23	1:A:191:TYR:N	2.32	0.45
1:A:41:TYR:CZ	1:A:137:GLY:HA2	2.52	0.45
1:A:58:GLN:C	1:A:60:PHE:H	2.20	0.45
1:A:149:TRP:CG	1:A:152:LYS:HG3	2.51	0.45
1:A:269:ARG:HD3	1:A:270:TYR:CE2	2.52	0.45
1:A:273:LEU:HA	1:A:276:SER:OG	2.17	0.45
1:A:176:PHE:CE1	1:A:190:LEU:N	2.85	0.45
1:A:183:PHE:CB	1:A:191:TYR:HH	2.29	0.44
1:A:398:VAL:O	1:A:402:ALA:HB3	2.17	0.44
1:A:42:TYR:CD1	1:A:165:HIS:HA	2.52	0.44
1:A:335:ALA:HB1	1:A:339:TYR:CE1	2.53	0.44
1:A:357:ILE:O	1:A:361:ILE:HG12	2.17	0.44
1:A:212:PRO:CG	1:A:213:GLN:N	2.80	0.44
1:A:53:PRO:HB3	1:A:278:THR:CB	2.48	0.44
1:A:368:ILE:O	1:A:371:HIS:N	2.50	0.44
1:A:45:ARG:NE	1:A:134:GLN:HG2	2.33	0.44
1:A:306:THR:HG23	1:A:307:LEU:N	2.31	0.44
1:A:427:ALA:O	1:A:431:LEU:HB2	2.18	0.44
1:A:26:LEU:HA	1:A:26:LEU:HD22	1.86	0.44
1:A:69:LEU:CD2	1:A:72:ILE:HD13	2.40	0.43
1:A:14:LEU:O	1:A:14:LEU:HG	2.18	0.43
1:A:339:TYR:CE2	1:A:349:VAL:HG12	2.53	0.43
1:A:339:TYR:CZ	1:A:353:CYS:HA	2.52	0.43
1:A:335:ALA:HB1	1:A:339:TYR:HE1	1.82	0.43
1:A:275:TRP:CH2	1:A:404:VAL:HB	2.53	0.43
1:A:44:VAL:O	1:A:47:ASN:ND2	2.51	0.43
1:A:218:PRO:O	1:A:220:ILE:HG12	2.17	0.43
1:A:114:PRO:O	1:A:115:TRP:C	2.56	0.43
1:A:269:ARG:HB2	1:A:361:ILE:CG2	2.46	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:346:ASN:HD21	1:A:350:ASP:CB	2.31	0.43
1:A:25:ARG:O	1:A:29:GLN:HB2	2.18	0.43
1:A:356:VAL:O	1:A:357:ILE:C	2.57	0.43
1:A:408:VAL:O	1:A:411:PHE:O	2.36	0.43
1:A:188:ALA:O	1:A:191:TYR:CE1	2.72	0.43
1:A:252:ASN:OD1	1:A:255:LEU:HB2	2.19	0.43
1:A:75:ALA:HB3	1:A:131:GLY:HA3	2.01	0.43
1:A:269:ARG:HG3	1:A:361:ILE:CD1	2.45	0.42
1:A:450:VAL:HG12	1:A:451:PRO:N	2.33	0.42
1:A:52:MET:CE	1:A:69:LEU:HD12	2.50	0.42
1:A:32:LEU:N	1:A:32:LEU:HD12	2.34	0.42
1:A:52:MET:HE2	1:A:69:LEU:HD12	2.01	0.42
1:A:30:ILE:HG23	1:A:31:PHE:N	2.34	0.42
1:A:251:PRO:O	1:A:252:ASN:C	2.58	0.42
1:A:241:ALA:HB3	1:A:243:GLN:HE21	1.84	0.42
1:A:16:ALA:O	1:A:17:ALA:HB3	2.19	0.42
1:A:100:LEU:O	1:A:103:ALA:N	2.53	0.42
1:A:50:LEU:HD11	1:A:173:PRO:HB3	2.01	0.42
1:A:303:ILE:HB	1:A:304:PRO:CD	2.39	0.42
1:A:142:GLY:O	1:A:146:VAL:HG23	2.20	0.42
1:A:136:MET:O	1:A:140:PRO:HD2	2.19	0.41
1:A:275:TRP:HA	1:A:278:THR:HG22	2.01	0.41
1:A:120:ILE:H	1:A:123:MET:HE2	1.84	0.41
1:A:15:PRO:HA	1:A:19:ILE:HG13	2.02	0.41
1:A:279:TYR:HD1	1:A:413:TRP:CE2	2.38	0.41
1:A:367:LEU:HA	1:A:367:LEU:HD12	1.58	0.41
1:A:176:PHE:CZ	1:A:189:ALA:HB3	2.54	0.41
1:A:436:ILE:HG22	1:A:437:GLY:N	2.35	0.41
1:A:79:SER:OG	1:A:80:LYS:N	2.53	0.41
1:A:335:ALA:HA	1:A:338:VAL:CG1	2.50	0.41
1:A:332:VAL:CG2	1:A:333:THR:N	2.84	0.41
1:A:328:PHE:CE2	1:A:360:LEU:HD22	2.55	0.41
1:A:426:LEU:O	1:A:427:ALA:C	2.58	0.41
1:A:122:VAL:HG13	1:A:123:MET:N	2.35	0.41
1:A:153:GLU:O	1:A:154:ARG:C	2.59	0.41
1:A:212:PRO:O	1:A:216:GLY:HA2	2.20	0.41
1:A:53:PRO:HB3	1:A:278:THR:HB	2.01	0.41
1:A:220:ILE:C	1:A:222:GLU:N	2.71	0.41
1:A:141:CYS:O	1:A:145:MET:HG3	2.21	0.41
1:A:388:THR:OG1	1:A:389:GLY:N	2.53	0.41
1:A:123:MET:O	1:A:124:PHE:C	2.57	0.41
1:A:397:SER:HA	1:A:400:ALA:HB3	2.01	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:108:LEU:CD2	1:A:112:PHE:HE1	2.33	0.41
1:A:56:VAL:HG21	1:A:61:SER:CB	2.50	0.41
1:A:140:PRO:C	1:A:142:GLY:N	2.72	0.41
1:A:328:PHE:O	1:A:332:VAL:CG1	2.64	0.41
1:A:85:SER:O	1:A:88:ASP:HB3	2.21	0.41
1:A:74:ILE:HG13	1:A:74:ILE:H	1.55	0.41
1:A:248:TYR:CD1	1:A:248:TYR:N	2.89	0.41
1:A:269:ARG:CB	1:A:361:ILE:HG21	2.49	0.40
1:A:252:ASN:HA	1:A:252:ASN:HD22	1.60	0.40
1:A:62:ARG:O	1:A:409:ASP:OD2	2.39	0.40
1:A:316:VAL:HG12	1:A:317:PHE:N	2.36	0.40
1:A:15:PRO:HA	1:A:19:ILE:CG1	2.51	0.40
1:A:355:ILE:O	1:A:359:PHE:HB2	2.21	0.40
1:A:88:ASP:O	1:A:143:ARG:NE	2.53	0.40
1:A:378:LYS:HG3	1:A:378:LYS:H	1.55	0.40
1:A:334:ILE:CG2	1:A:335:ALA:N	2.84	0.40
1:A:177:LEU:C	1:A:177:LEU:HD13	2.42	0.40
1:A:19:ILE:HG23	1:A:215:CYS:SG	2.62	0.40
1:A:211:THR:HA	1:A:213:GLN:NE2	2.36	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	430/451 (95%)	356 (83%)	50 (12%)	24 (6%)	3 25

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	184	ASN
1	A	211	THR
1	A	221	GLU
1	A	18	GLU

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Mol	Chain	Res	Type
1	A	59	GLY
1	A	100	LEU
1	A	251	PRO
1	A	253	LYS
1	A	9	PRO
1	A	212	PRO
1	A	242	LYS
1	A	243	GLN
1	A	15	PRO
1	A	137	GLY
1	A	344	ALA
1	A	346	ASN
1	A	378	LYS
1	A	120	ILE
1	A	114	PRO
1	A	158	VAL
1	A	56	VAL
1	A	219	PRO
1	A	396	GLY
1	A	7	PRO

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	345/360 (96%)	299 (87%)	46 (13%)	6 27

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	PHE
1	A	6	LYS
1	A	10	HIS
1	A	19	ILE
1	A	23	TYR
1	A	26	LEU
1	A	60	PHE
1	A	64	ASP

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Mol	Chain	Res	Type
1	A	69	LEU
1	A	79	SER
1	A	102	LEU
1	A	109	PHE
1	A	125	VAL
1	A	127	LEU
1	A	148	TRP
1	A	149	TRP
1	A	151	GLN
1	A	180	MET
1	A	186	TRP
1	A	187	HIS
1	A	190	LEU
1	A	195	PHE
1	A	196	CYS
1	A	203	PHE
1	A	210	ASP
1	A	252	ASN
1	A	262	ASN
1	A	266	TYR
1	A	272	ILE
1	A	279	TYR
1	A	295	TYR
1	A	299	GLU
1	A	327	PHE
1	A	329	MET
1	A	332	VAL
1	A	338	VAL
1	A	340	TRP
1	A	343	PRO
1	A	349	VAL
1	A	350	ASP
1	A	353	CYS
1	A	361	ILE
1	A	367	LEU
1	A	391	PHE
1	A	434	VAL
1	A	435	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	ASN
1	A	91	ASN
1	A	134	GLN
1	A	243	GLN
1	A	262	ASN
1	A	346	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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 ⓘ

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	434/451 (96%)	-0.24	2 (0%)	88 51	14, 74, 159, 200	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	222	GLU	2.2
1	A	223	TYR	2.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.