WORLDWIDE PDB
PROTEIN DATA BANK

Full wwPDB X-ray Structure Validation Report (i)

Jan 30, 2021 – 03:45 PM EST

PDB ID	:	3G79
Title	:	Crystal structure of NDP-N-acetyl-D-galactosaminuronic acid dehydrogenase from Methanosarcina mazei Go1
Authors	:	Malashkevich, V.N.; Toro, R.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on	:	2009-02-09
Resolution	:	2.40 Å(reported)

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

|		
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.16
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.16

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 2.40 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.

■ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution
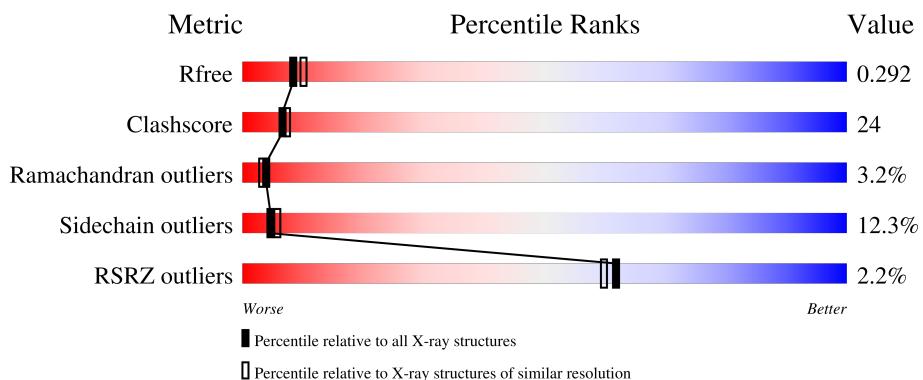
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 <=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	478	
1	B	478	
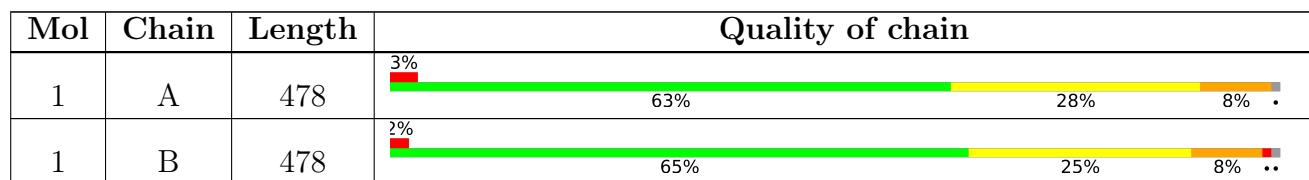
2 Entry composition ⓘ

There are 2 unique types of molecules in this entry. The entry contains 7421 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 NDP-N-acetyl-D-galactosaminuronic acid dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	475	Total 3621	C 2293	N 627	O 683	S 18	0	0	0
1	B	475	Total 3621	C 2293	N 627	O 683	S 18	0	0	0

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q8PXR2
A	2	SER	-	expression tag	UNP Q8PXR2
A	3	LEU	-	expression tag	UNP Q8PXR2
A	471	GLU	-	expression tag	UNP Q8PXR2
A	472	GLY	-	expression tag	UNP Q8PXR2
A	473	HIS	-	expression tag	UNP Q8PXR2
A	474	HIS	-	expression tag	UNP Q8PXR2
A	475	HIS	-	expression tag	UNP Q8PXR2
A	476	HIS	-	expression tag	UNP Q8PXR2
A	477	HIS	-	expression tag	UNP Q8PXR2
A	478	HIS	-	expression tag	UNP Q8PXR2
B	1	MET	-	expression tag	UNP Q8PXR2
B	2	SER	-	expression tag	UNP Q8PXR2
B	3	LEU	-	expression tag	UNP Q8PXR2
B	471	GLU	-	expression tag	UNP Q8PXR2
B	472	GLY	-	expression tag	UNP Q8PXR2
B	473	HIS	-	expression tag	UNP Q8PXR2
B	474	HIS	-	expression tag	UNP Q8PXR2
B	475	HIS	-	expression tag	UNP Q8PXR2
B	476	HIS	-	expression tag	UNP Q8PXR2
B	477	HIS	-	expression tag	UNP Q8PXR2
B	478	HIS	-	expression tag	UNP Q8PXR2

• Molecule 2 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	179	Total 179	O 179	0	0

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 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: NDP-N-acetyl-D-galactosaminuronic acid dehydrogenase

- Molecule 1: NDP-N-acetyl-D-galactosaminuronic acid dehydrogenase

S425 S428 D433 W434 V438 K441 A442 I446 I447 D448 G449 I453 K461 V464 Y465 K466 G467 I468 G469 R470 E471 G472 H473 H474 H478
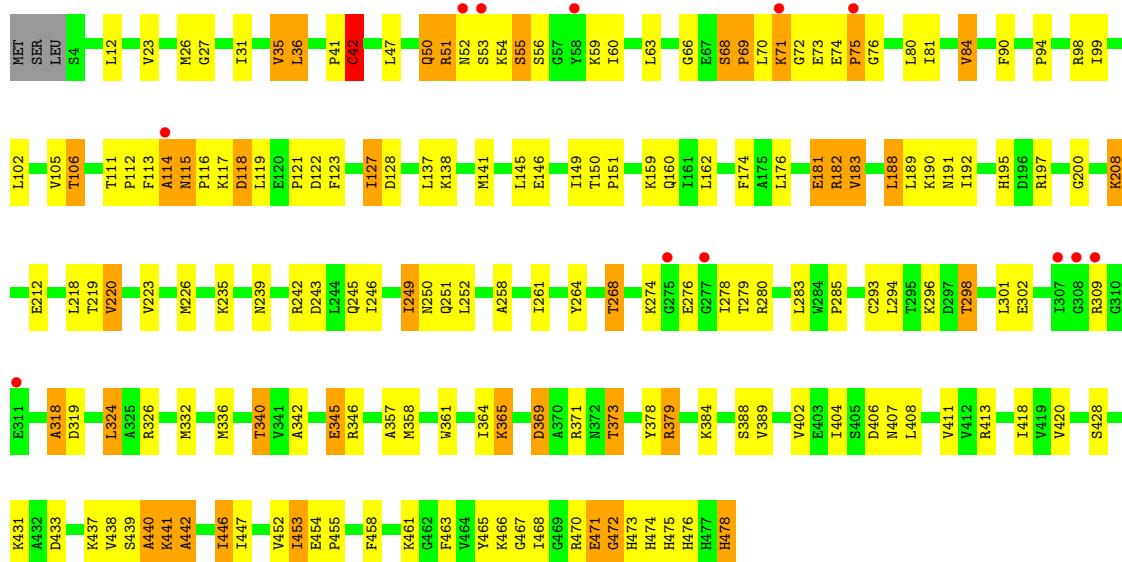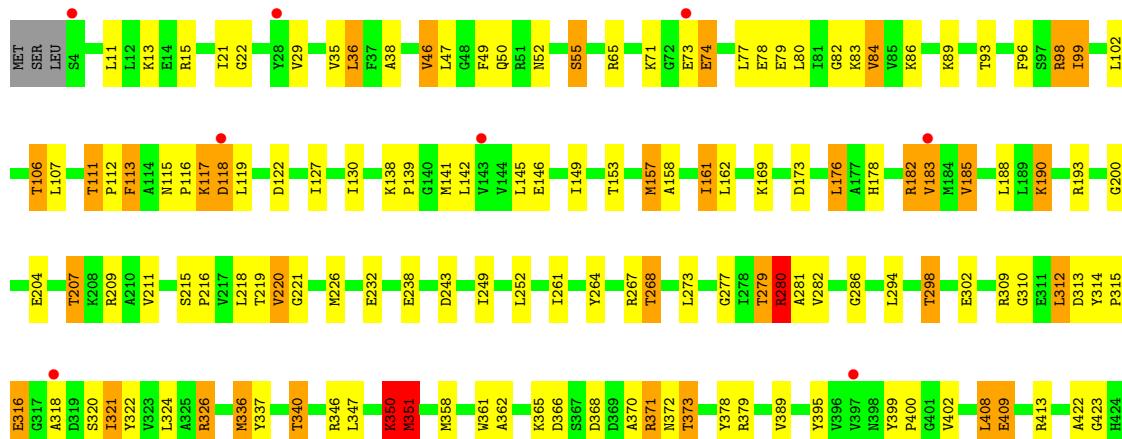
4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	78.83Å 88.72Å 151.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.71 – 2.40 19.71 – 2.40	Depositor EDS
% Data completeness (in resolution range)	95.4 (19.71-2.40) 95.4 (19.71-2.40)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.29 (at 2.41Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.222 , 0.296 0.218 , 0.292	Depositor DCC
R_{free} test set	2038 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å²)	44.2	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent k_{sol}(e/Å³), B_{sol}(Å²)	0.28 , 30.6	EDS
L-test for twinning[2]	$< \lvert L \rvert > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	7421	wwPDB-VP
Average B, all atoms (Å²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.64% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< \lvert L \rvert >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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	0.80	1/3690 (0.0%)	0.91	2/4986 (0.0%)
1	B	0.78	1/3690 (0.0%)	0.91	7/4986 (0.1%)
All	All	0.79	2/7380 (0.0%)	0.91	9/9972 (0.1%)

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

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	478	HIS	C-OXT	5.62	1.34	1.23
1	B	478	HIS	C-OXT	5.54	1.33	1.23

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	336	MET	CG-SD-CE	-7.98	87.43	100.20
1	B	326	ARG	NE-CZ-NH1	7.80	124.20	120.30
1	B	326	ARG	NE-CZ-NH2	-7.44	116.58	120.30
1	B	176	LEU	CA-CB-CG	6.18	129.51	115.30
1	B	280	ARG	NE-CZ-NH1	5.89	123.24	120.30
1	B	312	LEU	CA-CB-CG	5.55	128.06	115.30
1	B	98	ARG	N-CA-C	-5.19	96.99	111.00
1	A	379	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	A	98	ARG	NE-CZ-NH2	-5.07	117.77	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	112	PRO	Peptide
1	A	472	GLY	Peptide

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	3621	0	3655	205	0
1	B	3621	0	3655	160	0
2	A	179	0	0	10	0
All	All	7421	0	7310	343	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 24.

All (343) 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:471:GLU:CB	1:B:472:GLY:HA3	1.25	1.54
1:B:471:GLU:CB	1:B:472:GLY:CA	1.87	1.51
1:B:471:GLU:HB2	1:B:472:GLY:C	1.38	1.44
1:A:471:GLU:CB	1:A:472:GLY:CA	2.01	1.39
1:A:471:GLU:CB	1:A:472:GLY:HA3	1.42	1.35
1:B:471:GLU:HB3	1:B:472:GLY:CA	1.50	1.35
1:A:471:GLU:HB2	1:A:472:GLY:C	1.43	1.34
1:A:471:GLU:HB2	1:A:472:GLY:CA	1.62	1.29
1:A:342:ALA:O	1:A:345:GLU:HG3	1.31	1.27
1:B:336:MET:HE3	1:B:468:ILE:HD11	1.20	1.13
1:A:219:THR:O	1:A:220:VAL:HB	1.43	1.11
1:A:440:ALA:O	1:A:441:LYS:HB2	1.45	1.08
1:B:471:GLU:HB2	1:B:472:GLY:CA	1.63	1.08
1:A:74:GLU:O	1:A:76:GLY:N	1.88	1.07
1:A:342:ALA:O	1:A:345:GLU:CG	2.03	1.06
1:A:361:TRP:O	1:A:373:THR:HG21	1.55	1.06
1:B:361:TRP:O	1:B:373:THR:HG21	1.58	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:336:MET:HE3	1:A:468:ILE:HG13	1.36	1.03
1:A:114:ALA:HB1	1:A:115:ASN:HA	1.40	1.02
1:A:138:LYS:H	1:A:141:MET:HE2	1.20	1.02
1:B:115:ASN:C	1:B:117:LYS:H	1.62	1.02
2:A:643:HOH:O	1:B:313:ASP:HB2	1.61	1.00
1:B:115:ASN:O	1:B:117:LYS:N	1.94	0.99
1:B:471:GLU:CB	1:B:472:GLY:C	2.24	0.99
1:A:182:ARG:NE	1:B:273:LEU:HD21	1.79	0.98
1:B:336:MET:HE3	1:B:468:ILE:CD1	1.94	0.96
1:A:471:GLU:HB3	1:A:472:GLY:CA	1.73	0.96
1:A:113:PHE:HA	1:A:114:ALA:HB3	1.44	0.95
1:A:358:MET:HE1	1:A:379:ARG:HA	1.50	0.92
1:A:113:PHE:HB3	1:A:114:ALA:C	1.91	0.91
1:B:471:GLU:HB2	1:B:473:HIS:N	1.85	0.91
1:A:219:THR:OG1	1:A:219:THR:O	1.85	0.90
1:A:324:LEU:CD2	1:B:321:ILE:HG22	2.00	0.89
1:A:251:GLN:OE1	1:B:314:TYR:CE1	2.26	0.89
1:B:478:HIS:ND1	1:B:478:HIS:N	2.22	0.88
1:A:219:THR:O	1:A:220:VAL:CB	2.21	0.87
1:B:215:SER:HB2	1:B:216:PRO:HD3	1.57	0.87
1:A:342:ALA:HA	1:A:345:GLU:HG2	1.57	0.85
1:A:324:LEU:HD21	1:B:321:ILE:HG22	1.57	0.84
1:A:54:LYS:O	1:A:55:SER:HB3	1.76	0.84
1:A:138:LYS:H	1:A:141:MET:CE	1.90	0.84
1:A:441:LYS:O	1:A:442:ALA:CB	2.25	0.82
1:A:340:THR:HG21	1:A:378:TYR:OH	1.80	0.81
1:B:115:ASN:C	1:B:117:LYS:N	2.26	0.81
2:A:482:HOH:O	1:B:268:THR:HG22	1.79	0.81
1:A:467:GLY:O	1:A:474:HIS:CE1	2.34	0.81
1:A:471:GLU:HB2	1:A:473:HIS:N	1.95	0.81
1:A:471:GLU:CB	1:A:472:GLY:C	2.34	0.81
1:A:441:LYS:O	1:A:442:ALA:HB3	1.81	0.80
1:A:471:GLU:HB3	1:A:472:GLY:HA3	0.81	0.80
1:A:346:ARG:HH21	1:A:346:ARG:HG2	1.47	0.80
1:B:99:ILE:HA	1:B:102:LEU:HD12	1.63	0.80
1:A:235:LYS:NZ	1:A:293:CYS:SG	2.54	0.80
1:B:408:LEU:O	1:B:409:GLU:CB	2.27	0.79
1:A:137:LEU:HA	1:A:141:MET:CE	2.12	0.79
1:A:336:MET:CE	1:A:468:ILE:HG13	2.11	0.79
1:A:114:ALA:HB1	1:A:115:ASN:CA	2.14	0.78
1:A:358:MET:CE	1:A:379:ARG:HA	2.13	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:PHE:HB2	1:A:122:ASP:HB3	1.66	0.78
1:B:106:THR:HG21	1:B:146:GLU:OE1	1.84	0.78
1:B:408:LEU:O	1:B:409:GLU:HB3	1.81	0.78
1:B:340:THR:HG21	1:B:378:TYR:OH	1.85	0.77
1:B:366:ASP:HA	1:B:395:TYR:CE2	2.19	0.77
1:A:361:TRP:O	1:A:373:THR:CG2	2.32	0.76
1:B:336:MET:CE	1:B:468:ILE:CG1	2.63	0.76
1:B:157:MET:O	1:B:161:ILE:HG23	1.86	0.76
1:B:361:TRP:O	1:B:373:THR:CG2	2.33	0.75
1:B:264:TYR:O	1:B:268:THR:HG23	1.86	0.75
1:A:467:GLY:O	1:A:474:HIS:HE1	1.70	0.75
1:A:113:PHE:HB3	1:A:114:ALA:O	1.87	0.74
1:B:336:MET:HE1	1:B:468:ILE:CG1	2.17	0.74
1:B:200:GLY:HA2	1:B:226:MET:O	1.87	0.73
1:B:467:GLY:O	1:B:474:HIS:CE1	2.41	0.72
1:B:80:LEU:O	1:B:84:VAL:HG23	1.89	0.72
1:A:336:MET:HE3	1:A:468:ILE:CG1	2.17	0.72
1:A:251:GLN:OE1	1:B:314:TYR:HE1	1.71	0.72
1:B:471:GLU:OE1	1:B:471:GLU:N	2.22	0.72
1:A:84:VAL:HG13	1:A:90:PHE:HB3	1.71	0.72
1:B:182:ARG:NH2	1:B:238:GLU:OE1	2.23	0.72
1:A:440:ALA:O	1:A:441:LYS:CB	2.28	0.71
1:A:138:LYS:N	1:A:141:MET:HE2	2.02	0.71
1:A:466:LYS:HA	1:A:474:HIS:CD2	2.26	0.71
1:B:112:PRO:O	1:B:113:PHE:HB2	1.89	0.71
1:A:324:LEU:HD22	1:B:324:LEU:HD23	1.73	0.71
1:B:79:GLU:OE1	1:B:79:GLU:N	2.15	0.71
1:A:114:ALA:CB	1:A:115:ASN:HA	2.20	0.70
1:A:243:ASP:OD2	1:A:326:ARG:HD2	1.92	0.70
1:B:38:ALA:CB	1:B:84:VAL:HG11	2.22	0.70
1:A:113:PHE:CA	1:A:114:ALA:HB3	2.20	0.70
1:B:261:ILE:C	1:B:471:GLU:HA	2.12	0.70
1:A:113:PHE:O	1:A:121:PRO:HA	1.92	0.69
1:A:250:ASN:HB3	1:A:332:MET:CE	2.23	0.69
1:A:113:PHE:CD1	1:A:114:ALA:HB3	2.27	0.69
1:A:50:GLN:NE2	1:A:51:ARG:O	2.26	0.69
1:B:38:ALA:HB3	1:B:84:VAL:HG11	1.75	0.68
1:A:245:GLN:HG3	1:A:283:LEU:CD1	2.24	0.68
1:A:183:VAL:HG13	1:A:188:LEU:HD12	1.75	0.68
1:B:467:GLY:N	1:B:474:HIS:NE2	2.32	0.67
1:A:54:LYS:O	1:A:55:SER:CB	2.43	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:471:GLU:N	1:A:471:GLU:OE1	2.28	0.67
1:B:340:THR:HB	1:B:447:ILE:HG13	1.77	0.67
1:B:111:THR:HG23	1:B:111:THR:O	1.95	0.66
1:A:245:GLN:HG3	1:A:283:LEU:HD12	1.77	0.66
1:A:319:ASP:O	1:B:324:LEU:HD21	1.96	0.66
1:B:219:THR:OG1	1:B:219:THR:O	2.11	0.65
1:B:336:MET:HE1	1:B:468:ILE:HG12	1.78	0.65
1:A:218:LEU:CD1	1:A:223:VAL:HG23	2.26	0.65
1:A:261:ILE:C	1:A:471:GLU:HA	2.17	0.65
1:B:358:MET:CE	1:B:379:ARG:HA	2.27	0.65
1:A:35:VAL:CG1	1:A:80:LEU:HG	2.26	0.65
1:A:26:MET:HE1	1:A:90:PHE:HZ	1.62	0.65
1:A:102:LEU:HD12	1:A:105:VAL:HG22	1.79	0.65
1:A:470:ARG:HB2	1:A:474:HIS:NE2	2.13	0.64
1:B:169:LYS:N	1:B:173:ASP:OD2	2.21	0.64
1:A:35:VAL:HG13	1:A:80:LEU:HG	1.78	0.64
1:A:50:GLN:O	1:A:94:PRO:HA	1.97	0.64
1:B:336:MET:HE3	1:B:468:ILE:CG1	2.26	0.64
1:A:242:ARG:HG2	1:A:294:LEU:HD21	1.78	0.64
1:A:358:MET:HE3	1:A:389:VAL:HG11	1.79	0.64
1:A:66:GLY:HA2	1:A:81:ILE:HG22	1.80	0.64
1:B:35:VAL:CG1	1:B:80:LEU:HG	2.27	0.64
1:B:106:THR:CG2	1:B:146:GLU:OE1	2.46	0.63
1:A:324:LEU:HD21	1:B:321:ILE:HA	1.79	0.63
1:A:200:GLY:HA2	1:A:226:MET:O	1.99	0.63
1:A:324:LEU:HD21	1:B:321:ILE:CG2	2.28	0.62
1:A:137:LEU:HD12	1:A:141:MET:HE2	1.81	0.62
1:A:113:PHE:HA	1:A:114:ALA:CB	2.25	0.62
1:A:342:ALA:HA	1:A:345:GLU:CG	2.30	0.62
1:B:470:ARG:HB2	1:B:474:HIS:NE2	2.15	0.62
1:B:107:LEU:HD13	1:B:145:LEU:HD23	1.81	0.61
1:B:185:VAL:HA	1:B:188:LEU:HD13	1.80	0.61
1:B:336:MET:O	1:B:340:THR:HG22	2.01	0.61
1:B:314:TYR:CG	1:B:315:PRO:HD2	2.36	0.60
1:A:365:LYS:NZ	2:A:514:HOH:O	2.32	0.60
1:B:358:MET:HE1	1:B:379:ARG:HA	1.83	0.60
1:B:157:MET:HG3	1:B:158:ALA:N	2.17	0.60
1:A:26:MET:CE	1:A:90:PHE:HZ	2.15	0.59
1:A:70:LEU:O	1:A:71:LYS:HB2	2.02	0.59
1:A:52:ASN:HB2	2:A:636:HOH:O	2.02	0.59
1:A:123:PHE:O	1:A:127:ILE:HG23	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:LEU:HD13	1:A:80:LEU:HD21	1.84	0.59
1:A:467:GLY:N	1:A:474:HIS:NE2	2.37	0.59
1:B:130:ILE:HB	1:B:161:ILE:HD11	1.85	0.59
1:B:466:LYS:HA	1:B:474:HIS:CD2	2.38	0.58
1:A:84:VAL:HG13	1:A:90:PHE:CB	2.33	0.58
1:A:218:LEU:HD11	1:A:223:VAL:HG23	1.84	0.58
1:A:114:ALA:HA	1:A:117:LYS:H	1.69	0.58
1:A:182:ARG:CZ	1:B:273:LEU:HD21	2.33	0.58
1:A:298:THR:O	1:A:302:GLU:HG2	2.03	0.58
1:B:99:ILE:HA	1:B:102:LEU:CD1	2.33	0.57
1:B:279:THR:O	1:B:280:ARG:HB2	2.04	0.57
1:A:111:THR:HG23	1:A:111:THR:O	2.03	0.57
1:A:68:SER:HB2	1:A:69:PRO:HD3	1.85	0.57
1:B:111:THR:CG2	1:B:111:THR:O	2.51	0.57
1:B:207:THR:O	1:B:211:VAL:HG23	2.05	0.57
1:A:73:GLU:O	1:A:74:GLU:C	2.43	0.57
1:A:318:ALA:O	1:A:319:ASP:HB2	2.04	0.57
1:B:11:LEU:O	1:B:15:ARG:HG2	2.06	0.56
1:B:350:LYS:O	1:B:351:MET:HB2	2.05	0.56
1:B:347:LEU:HD21	1:B:464:VAL:HG21	1.87	0.56
1:A:182:ARG:HE	1:B:273:LEU:HD21	1.66	0.56
1:A:208:LYS:O	1:A:212:GLU:HG3	2.06	0.56
1:A:251:GLN:HE22	1:B:320:SER:HB2	1.71	0.55
1:A:74:GLU:HG3	1:A:189:LEU:H	1.71	0.55
1:B:336:MET:CE	1:B:468:ILE:HG12	2.33	0.55
1:A:26:MET:HE1	1:A:90:PHE:CZ	2.41	0.55
1:B:441:LYS:O	1:B:442:ALA:CB	2.54	0.55
1:A:250:ASN:HB3	1:A:332:MET:HE3	1.87	0.55
1:A:358:MET:CE	1:A:389:VAL:HG11	2.36	0.55
1:A:439:SER:C	1:A:440:ALA:O	2.44	0.55
1:A:342:ALA:C	1:A:345:GLU:CG	2.74	0.54
1:B:219:THR:O	1:B:220:VAL:CB	2.55	0.54
1:A:41:PRO:O	1:A:42:CYS:HB3	2.06	0.54
1:A:470:ARG:O	1:A:471:GLU:O	2.25	0.54
1:A:84:VAL:CG1	1:A:90:PHE:HB3	2.38	0.54
1:B:336:MET:HE1	1:B:468:ILE:HG13	1.89	0.54
1:A:251:GLN:HG2	1:B:322:TYR:CZ	2.42	0.54
1:A:68:SER:HB2	1:A:69:PRO:CD	2.37	0.54
1:A:56:SER:O	1:A:59:LYS:HG3	2.08	0.54
1:A:68:SER:CB	1:A:69:PRO:CD	2.86	0.54
1:B:219:THR:O	1:B:220:VAL:HB	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:336:MET:HE2	1:B:336:MET:HA	1.88	0.54
1:B:98:ARG:O	1:B:99:ILE:HG13	2.07	0.54
1:A:250:ASN:HB3	1:A:332:MET:HE1	1.89	0.53
1:B:215:SER:HB2	1:B:216:PRO:CD	2.35	0.53
1:B:467:GLY:O	1:B:474:HIS:HE1	1.88	0.53
1:A:74:GLU:C	1:A:76:GLY:H	2.00	0.53
1:A:324:LEU:HD22	1:B:324:LEU:CD2	2.38	0.53
2:A:585:HOH:O	1:B:55:SER:HB2	2.08	0.53
1:A:116:PRO:O	1:A:117:LYS:HB2	2.09	0.53
1:A:358:MET:HE3	1:A:379:ARG:HG3	1.90	0.52
1:A:137:LEU:HD12	1:A:141:MET:CE	2.39	0.52
1:A:138:LYS:N	1:A:141:MET:CE	2.64	0.52
1:A:342:ALA:O	1:A:345:GLU:CD	2.48	0.52
1:A:418:ILE:CG2	1:A:446:ILE:HD12	2.40	0.52
1:A:258:ALA:O	1:B:310:GLY:HA3	2.10	0.51
1:A:342:ALA:C	1:A:345:GLU:HG3	2.21	0.51
1:A:41:PRO:O	1:A:42:CYS:CB	2.58	0.51
1:A:324:LEU:CD2	1:B:321:ILE:CG2	2.82	0.51
1:A:182:ARG:CZ	1:B:273:LEU:CD2	2.89	0.51
1:A:111:THR:HG22	2:A:645:HOH:O	2.11	0.51
1:A:191:ASN:O	1:A:195:HIS:HB2	2.10	0.51
1:B:358:MET:CE	1:B:389:VAL:HG11	2.41	0.51
1:B:471:GLU:HB3	1:B:472:GLY:HA3	0.54	0.51
1:B:243:ASP:OD2	1:B:326:ARG:HD2	2.11	0.51
1:A:145:LEU:HD12	1:A:149:ILE:HD13	1.93	0.51
1:A:433:ASP:OD1	1:A:461:LYS:NZ	2.38	0.51
1:A:336:MET:O	1:A:340:THR:HG22	2.11	0.50
1:B:470:ARG:O	1:B:471:GLU:C	2.50	0.50
1:A:470:ARG:O	1:A:471:GLU:C	2.49	0.50
1:B:337:TYR:HA	1:B:340:THR:HG23	1.93	0.50
1:B:98:ARG:O	1:B:99:ILE:CB	2.59	0.50
1:A:113:PHE:HB2	1:A:122:ASP:CB	2.40	0.50
1:A:31:ILE:O	1:A:35:VAL:HB	2.11	0.50
1:B:358:MET:HE3	1:B:389:VAL:HG11	1.94	0.50
1:B:38:ALA:HB1	1:B:84:VAL:HG11	1.94	0.50
1:A:264:TYR:O	1:A:268:THR:CG2	2.59	0.50
1:A:53:SER:O	1:A:54:LYS:C	2.51	0.49
1:A:127:ILE:HG13	1:A:128:ASP:N	2.26	0.49
1:B:358:MET:HE3	1:B:379:ARG:HA	1.93	0.49
1:A:116:PRO:C	1:A:117:LYS:HD2	2.32	0.49
1:A:26:MET:HE2	1:A:63:LEU:HD13	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:118:ASP:N	1:B:118:ASP:OD2	2.45	0.49
1:B:409:GLU:HB2	1:B:434:TRP:HZ2	1.78	0.49
1:A:23:VAL:HG13	1:A:106:THR:HG22	1.94	0.49
1:A:342:ALA:CA	1:A:345:GLU:HG2	2.33	0.49
1:A:402:VAL:HG12	1:A:404:ILE:CD1	2.43	0.49
1:A:324:LEU:HD23	1:B:321:ILE:HG22	1.91	0.49
1:B:36:LEU:HD13	1:B:80:LEU:HD21	1.93	0.49
1:A:418:ILE:HG22	1:A:446:ILE:HD12	1.95	0.49
1:B:82:GLY:O	1:B:86:LYS:HG3	2.13	0.48
1:A:71:LYS:HG2	2:A:610:HOH:O	2.12	0.48
1:A:138:LYS:O	1:A:141:MET:HG3	2.14	0.48
1:B:346:ARG:NH1	1:B:478:HIS:CE1	2.82	0.48
1:B:22:GLY:HA2	1:B:47:LEU:O	2.13	0.48
1:B:470:ARG:O	1:B:471:GLU:O	2.32	0.47
1:A:102:LEU:CD1	1:A:105:VAL:HG22	2.44	0.47
1:A:47:LEU:HD12	1:A:102:LEU:HD21	1.95	0.47
1:A:433:ASP:HA	1:A:461:LYS:HE2	1.95	0.47
1:B:78:GLU:N	1:B:79:GLU:OE1	2.47	0.47
1:A:239:ASN:ND2	1:A:293:CYS:HB3	2.29	0.47
1:A:418:ILE:O	1:A:446:ILE:HA	2.15	0.47
1:A:251:GLN:NE2	1:B:320:SER:HB2	2.29	0.47
1:A:99:ILE:HG23	1:A:105:VAL:HG21	1.96	0.46
1:B:117:LYS:O	1:B:118:ASP:C	2.53	0.46
1:B:112:PRO:O	1:B:113:PHE:CB	2.60	0.46
1:A:26:MET:CE	1:A:63:LEU:HD13	2.45	0.46
1:A:53:SER:OG	2:A:609:HOH:O	2.19	0.46
1:A:113:PHE:HB3	1:A:114:ALA:CA	2.46	0.46
1:A:118:ASP:HB2	1:A:119:LEU:HA	1.97	0.46
1:B:13:LYS:HA	1:B:13:LYS:HD3	1.73	0.46
1:B:96:PHE:O	1:B:98:ARG:O	2.33	0.46
1:A:113:PHE:HD1	1:A:114:ALA:HB3	1.75	0.46
1:A:264:TYR:O	1:A:268:THR:HG23	2.16	0.46
1:B:314:TYR:CD1	1:B:315:PRO:HD2	2.51	0.46
1:B:365:LYS:HD3	1:B:423:GLY:O	2.15	0.46
1:A:246:ILE:O	1:A:249:ILE:HD12	2.16	0.45
1:B:74:GLU:OE1	1:B:188:LEU:HB2	2.16	0.45
1:A:150:THR:O	1:A:151:PRO:C	2.51	0.45
1:A:137:LEU:HD21	1:A:174:PHE:CD2	2.51	0.45
1:B:346:ARG:NH1	1:B:478:HIS:HE1	2.13	0.45
1:B:79:GLU:H	1:B:79:GLU:CD	2.05	0.45
1:A:245:GLN:HG3	1:A:283:LEU:HD11	1.95	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:447:ILE:HG13	1:A:466:LYS:HB2	1.97	0.45
1:B:29:VAL:HG22	1:B:183:VAL:HG11	1.99	0.45
1:A:27:GLY:O	1:A:31:ILE:HG13	2.17	0.45
1:B:107:LEU:HD13	1:B:145:LEU:CD2	2.46	0.45
1:B:219:THR:O	1:B:220:VAL:HG23	2.17	0.45
1:B:261:ILE:O	1:B:471:GLU:HA	2.16	0.45
1:B:409:GLU:HB2	1:B:434:TRP:CZ2	2.52	0.45
1:B:158:ALA:O	1:B:162:LEU:HB2	2.17	0.45
1:B:190:LYS:HA	1:B:193:ARG:HB2	1.98	0.45
1:B:413:ARG:HA	1:B:438:VAL:O	2.18	0.44
1:B:138:LYS:HB3	1:B:139:PRO:HD2	1.99	0.44
1:B:49:PHE:HA	1:B:93:THR:O	2.18	0.44
1:A:342:ALA:CA	1:A:345:GLU:CG	2.94	0.44
1:A:246:ILE:HA	1:A:249:ILE:HD11	1.99	0.44
1:A:118:ASP:HA	1:A:296:LYS:NZ	2.33	0.44
1:A:137:LEU:HA	1:A:141:MET:HE2	1.98	0.44
1:B:279:THR:HG22	1:B:281:ALA:H	1.81	0.44
1:B:464:VAL:HG22	1:B:478:HIS:CD2	2.52	0.44
1:A:36:LEU:HD13	1:A:80:LEU:CD2	2.46	0.43
1:A:358:MET:CE	1:A:379:ARG:CA	2.93	0.43
1:A:420:VAL:HG21	1:A:453:ILE:CD1	2.48	0.43
1:B:38:ALA:O	1:B:89:LYS:HE2	2.18	0.43
1:A:452:VAL:HG23	1:A:453:ILE:HD12	2.00	0.43
1:A:115:ASN:O	1:A:115:ASN:ND2	2.49	0.43
1:A:242:ARG:NH2	1:B:277:GLY:O	2.52	0.43
1:B:446:ILE:HD11	1:B:453:ILE:HD13	2.00	0.43
2:A:533:HOH:O	1:B:478:HIS:HA	2.19	0.43
1:A:106:THR:CG2	1:A:146:GLU:OE1	2.66	0.43
1:A:357:ALA:HB1	1:A:411:VAL:HG12	2.01	0.43
1:A:369:ASP:HB2	1:A:371:ARG:HH11	1.84	0.43
1:B:399:TYR:HA	1:B:400:PRO:HD3	1.92	0.43
1:B:286:GLY:HA2	1:B:449:GLY:O	2.19	0.43
1:B:80:LEU:HA	1:B:83:LYS:HG3	1.99	0.43
1:A:261:ILE:O	1:A:471:GLU:HA	2.19	0.43
1:A:60:ILE:HD13	1:A:94:PRO:HD3	2.01	0.42
1:B:71:LYS:HB3	1:B:71:LYS:HE3	1.80	0.42
1:A:346:ARG:NH2	1:A:346:ARG:HG2	2.24	0.42
1:A:181:GLU:HG2	1:A:197:ARG:NE	2.34	0.42
1:B:218:LEU:HD13	1:B:221:GLY:C	2.40	0.42
1:B:21:ILE:O	1:B:46:VAL:HA	2.19	0.42
1:B:361:TRP:CD1	1:B:370:ALA:CB	3.03	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:ARG:O	1:B:99:ILE:CG1	2.68	0.42
1:A:52:ASN:CG	1:A:52:ASN:O	2.57	0.42
1:A:69:PRO:HB2	1:A:70:LEU:H	1.64	0.42
1:A:183:VAL:CG1	1:A:188:LEU:HD12	2.48	0.42
1:A:433:ASP:OD1	1:A:461:LYS:CE	2.68	0.42
1:B:362:ALA:O	1:B:422:ALA:HB2	2.20	0.42
1:B:178:HIS:CD2	1:B:232:GLU:HG3	2.55	0.41
1:A:358:MET:HE1	1:A:379:ARG:CA	2.36	0.41
1:A:476:HIS:CE1	1:A:478:HIS:HB3	2.55	0.41
1:B:115:ASN:H	1:B:117:LYS:HB2	1.85	0.41
1:B:138:LYS:HB2	1:B:141:MET:CE	2.50	0.41
1:B:182:ARG:CZ	1:B:238:GLU:OE1	2.67	0.41
1:A:74:GLU:N	1:A:75:PRO:HD3	2.35	0.41
1:A:250:ASN:OD1	1:A:285:PRO:HA	2.21	0.41
1:A:264:TYR:O	1:A:268:THR:HG22	2.20	0.41
1:A:402:VAL:HG12	1:A:404:ILE:HD11	2.02	0.41
1:B:267:ARG:HG3	1:B:282:VAL:HG11	2.03	0.41
1:B:98:ARG:O	1:B:99:ILE:HB	2.19	0.41
1:A:74:GLU:OE2	1:A:190:LYS:HE3	2.20	0.41
1:A:245:GLN:OE1	1:B:280:ARG:HG3	2.21	0.41
1:A:324:LEU:HG	1:B:321:ILE:HG23	2.02	0.41
1:B:273:LEU:O	1:B:279:THR:HA	2.20	0.41
1:A:420:VAL:HG21	1:A:453:ILE:HD13	2.01	0.41
1:A:458:PHE:O	1:A:463:PHE:HB2	2.21	0.41
1:A:465:TYR:OH	1:A:470:ARG:HG3	2.21	0.41
1:B:142:LEU:HD22	1:B:209:ARG:HG3	2.03	0.41
1:B:145:LEU:HD12	1:B:149:ILE:HD13	2.02	0.41
2:A:587:HOH:O	1:B:294:LEU:O	2.22	0.41
1:A:378:TYR:C	1:A:378:TYR:CD2	2.94	0.41
1:A:454:GLU:HA	1:A:455:PRO:HD3	1.96	0.41
1:B:371:ARG:O	1:B:372:ASN:HB2	2.20	0.41
1:A:52:ASN:C	1:A:52:ASN:ND2	2.74	0.40
1:B:149:ILE:HD12	1:B:153:THR:HB	2.02	0.40
1:A:190:LYS:HE2	1:A:190:LYS:HB3	1.62	0.40
1:A:73:GLU:O	1:A:75:PRO:N	2.54	0.40
1:B:298:THR:O	1:B:302:GLU:HG3	2.21	0.40
1:A:181:GLU:OE2	1:A:192:ILE:HA	2.20	0.40
1:A:413:ARG:HA	1:A:438:VAL:HG12	2.02	0.40
1:B:433:ASP:OD1	1:B:461:LYS:HE3	2.21	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	473/478 (99%)	434 (92%)	23 (5%)	16 (3%)	3	3
1	B	473/478 (99%)	435 (92%)	24 (5%)	14 (3%)	4	3
All	All	946/956 (99%)	869 (92%)	47 (5%)	30 (3%)	4	3

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	68	SER
1	A	69	PRO
1	A	71	LYS
1	A	75	PRO
1	A	220	VAL
1	A	441	LYS
1	A	442	ALA
1	A	471	GLU
1	B	99	ILE
1	B	220	VAL
1	B	351	MET
1	B	409	GLU
1	B	471	GLU
1	A	42	CYS
1	A	55	SER
1	A	72	GLY
1	A	114	ALA
1	A	118	ASP
1	B	117	LYS
1	B	318	ALA
1	B	442	ALA
1	A	318	ALA
1	A	365	LYS
1	B	316	GLU
1	A	440	ALA

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Mol	Chain	Res	Type
1	B	113	PHE
1	B	116	PRO
1	B	350	LYS
1	B	280	ARG
1	B	74	GLU

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	383/387 (99%)	336 (88%)	47 (12%)	4	6
1	B	383/387 (99%)	336 (88%)	47 (12%)	4	6
All	All	766/774 (99%)	672 (88%)	94 (12%)	4	6

All (94) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LEU
1	A	35	VAL
1	A	36	LEU
1	A	42	CYS
1	A	50	GLN
1	A	51	ARG
1	A	84	VAL
1	A	106	THR
1	A	115	ASN
1	A	127	ILE
1	A	159	LYS
1	A	160	GLN
1	A	162	LEU
1	A	176	LEU
1	A	181	GLU
1	A	182	ARG
1	A	183	VAL
1	A	188	LEU

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Mol	Chain	Res	Type
1	A	208	LYS
1	A	249	ILE
1	A	252	LEU
1	A	268	THR
1	A	274	LYS
1	A	276	GLU
1	A	278	ILE
1	A	279	THR
1	A	280	ARG
1	A	298	THR
1	A	301	LEU
1	A	309	ARG
1	A	324	LEU
1	A	340	THR
1	A	345	GLU
1	A	364	ILE
1	A	369	ASP
1	A	373	THR
1	A	384	LYS
1	A	388	SER
1	A	406	ASP
1	A	407	ASN
1	A	408	LEU
1	A	428	SER
1	A	431	LYS
1	A	437	LYS
1	A	446	ILE
1	A	453	ILE
1	A	475	HIS
1	B	36	LEU
1	B	46	VAL
1	B	50	GLN
1	B	52	ASN
1	B	55	SER
1	B	65	ARG
1	B	73	GLU
1	B	77	LEU
1	B	84	VAL
1	B	106	THR
1	B	111	THR
1	B	118	ASP
1	B	119	LEU

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Mol	Chain	Res	Type
1	B	122	ASP
1	B	127	ILE
1	B	157	MET
1	B	161	ILE
1	B	176	LEU
1	B	182	ARG
1	B	183	VAL
1	B	185	VAL
1	B	190	LYS
1	B	204	GLU
1	B	207	THR
1	B	249	ILE
1	B	252	LEU
1	B	268	THR
1	B	279	THR
1	B	280	ARG
1	B	298	THR
1	B	309	ARG
1	B	312	LEU
1	B	316	GLU
1	B	321	ILE
1	B	340	THR
1	B	350	LYS
1	B	351	MET
1	B	368	ASP
1	B	371	ARG
1	B	373	THR
1	B	402	VAL
1	B	408	LEU
1	B	425	SER
1	B	428	SER
1	B	433	ASP
1	B	446	ILE
1	B	478	HIS

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

Mol	Chain	Res	Type
1	A	52	ASN
1	A	132	ASN
1	A	160	GLN
1	A	195	HIS

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Mol	Chain	Res	Type
1	A	251	GLN
1	A	398	ASN
1	B	50	GLN
1	B	52	ASN
1	B	191	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 monosaccharides 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	475/478 (99%)	-0.26	12 (2%)	57	55	21, 34, 60, 78	0
1	B	475/478 (99%)	-0.26	9 (1%)	66	64	20, 35, 61, 79	0
All	All	950/956 (99%)	-0.26	21 (2%)	62	60	20, 34, 61, 79	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	75	PRO	6.0
1	B	28	TYR	4.1
1	A	309	ARG	3.9
1	A	52	ASN	2.9
1	A	275	GLY	2.9
1	A	58	TYR	2.9
1	B	473	HIS	2.7
1	B	118	ASP	2.6
1	A	307	ILE	2.5
1	B	4	SER	2.5
1	B	183	VAL	2.5
1	A	308	GLY	2.4
1	A	311	GLU	2.4
1	A	114	ALA	2.4
1	B	397	VAL	2.4
1	A	277	GLY	2.3
1	B	73	GLU	2.3
1	A	71	LYS	2.3
1	B	143	VAL	2.1
1	B	318	ALA	2.0
1	A	53	SER	2.0

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 monosaccharides 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.