



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:43 PM GMT

PDB ID : 3RZF
Title : Crystal Structure of Inhibitor of kappaB kinase beta (I4122)
Authors : Xu, G.; Lo, Y.C.; Li, Q.; Napolitano, G.; Wu, X.; Jiang, X.; Dreano, M.; Karin, M.; Wu, H.
Deposited on : 2011-05-11
Resolution : 4.00 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

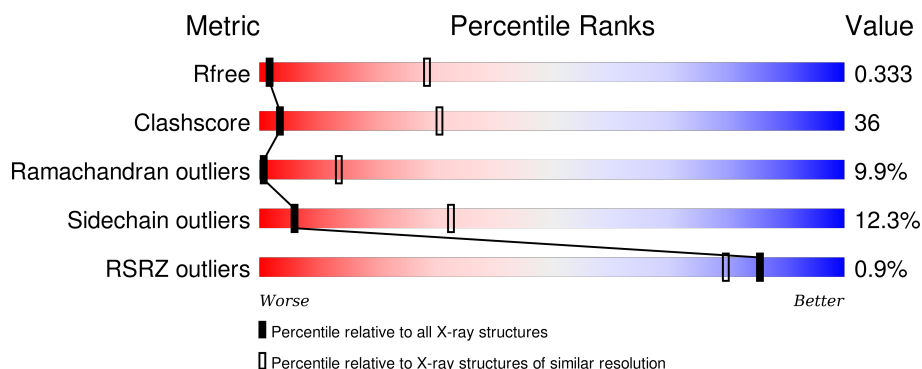
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.00 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}	91344	1010 (4.42-3.56)
Clashscore	102246	1052 (4.40-3.60)
Ramachandran outliers	100387	1005 (4.40-3.60)
Sidechain outliers	100360	1013 (4.42-3.58)
RSRZ outliers	91569	1013 (4.42-3.56)

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. 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	677	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	XNM	A	676	-	-	X	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4399 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.

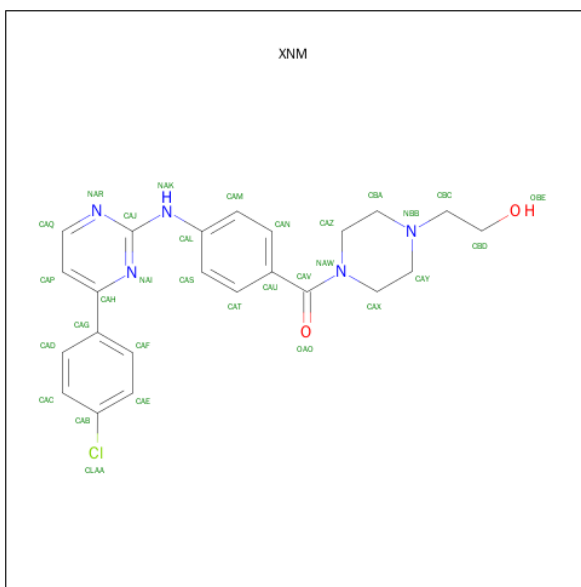
- Molecule 1 is a protein called MGC80376 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	541	Total	C	N	O	S	0	0	0
			4368	2764	779	799	26			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLN	-	EXPRESSION TAG	UNP Q6INT1
A	0	GLY	-	EXPRESSION TAG	UNP Q6INT1
A	1	GLY	-	EXPRESSION TAG	UNP Q6INT1
A	2	GLY	-	EXPRESSION TAG	UNP Q6INT1
A	3	ARG	-	EXPRESSION TAG	UNP Q6INT1
A	4	SER	-	EXPRESSION TAG	UNP Q6INT1
A	5	PRO	-	EXPRESSION TAG	UNP Q6INT1
A	6	SER	-	EXPRESSION TAG	UNP Q6INT1
A	7	LEU	-	EXPRESSION TAG	UNP Q6INT1
A	8	PRO	-	EXPRESSION TAG	UNP Q6INT1
A	9	THR	-	EXPRESSION TAG	UNP Q6INT1
A	10	GLN	-	EXPRESSION TAG	UNP Q6INT1
A	11	THR	-	EXPRESSION TAG	UNP Q6INT1
A	12	CYS	-	EXPRESSION TAG	UNP Q6INT1
A	13	GLY	-	EXPRESSION TAG	UNP Q6INT1
A	14	PRO	-	EXPRESSION TAG	UNP Q6INT1
A	15	TRP	-	EXPRESSION TAG	UNP Q6INT1
A	16	GLU	-	EXPRESSION TAG	UNP Q6INT1
A	177	GLU	SER	ENGINEERED MUTATION	UNP Q6INT1
A	181	GLU	SER	ENGINEERED MUTATION	UNP Q6INT1

- Molecule 2 is (4-{[4-(4-CHLOROPHENYL)PYRIMIDIN-2-YL]AMINO}PHENYL)[4-(2-H YDROXYETHYL)PIPERAZIN-1-YL]METHANONE (three-letter code: XNM) (formula: C₂₃H₂₄ClN₅O₂).

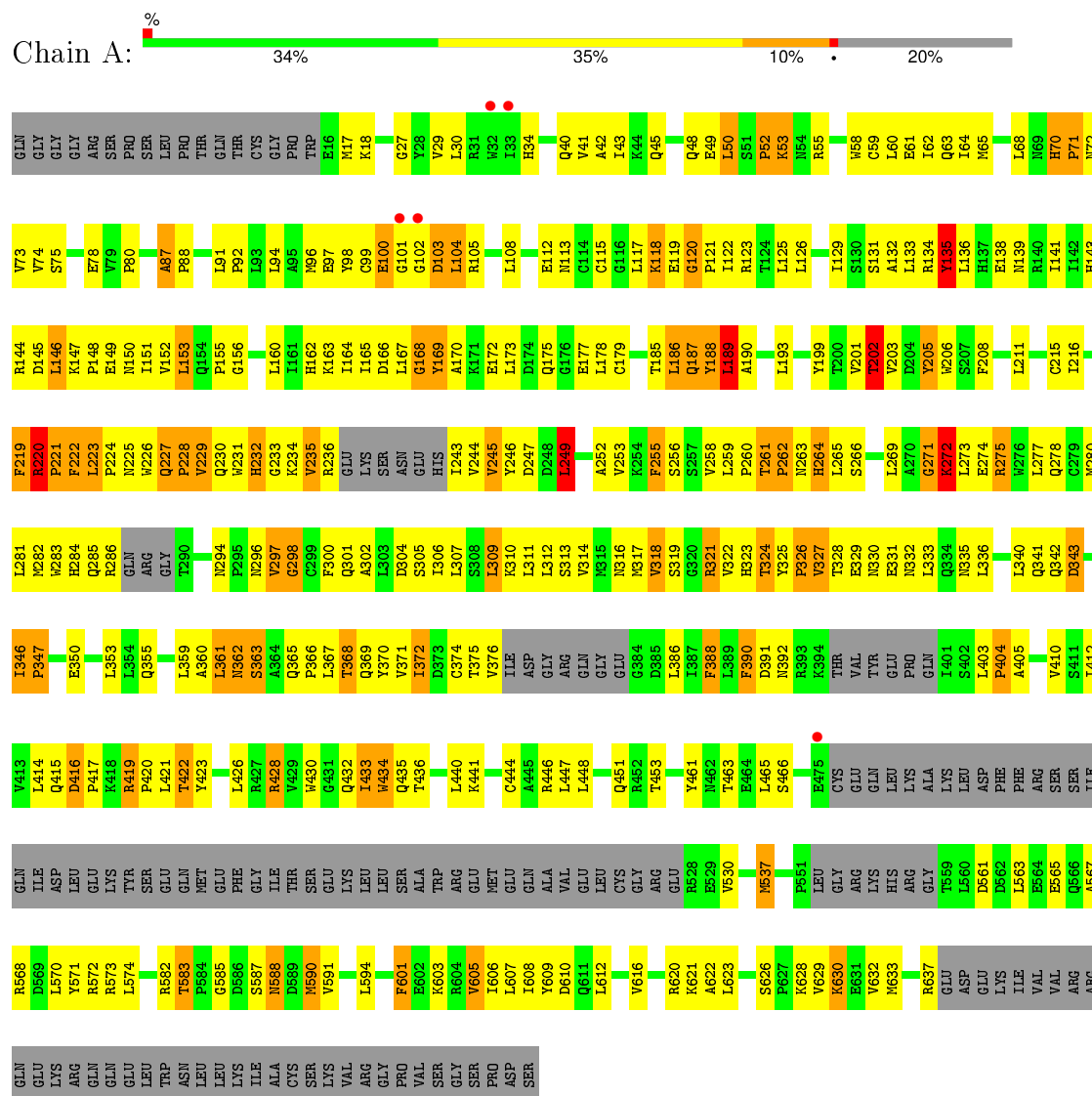


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	0	0
			31	23	1	5	2		

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: MGC80376 protein



4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	164.66 Å 164.66 Å 273.46 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.96 – 4.00 36.58 – 3.90	Depositor EDS
% Data completeness (in resolution range)	83.7 (14.96-4.00) 80.2 (36.58-3.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.52 (at 3.87 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.270 , 0.340 0.262 , 0.333	Depositor DCC
R_{free} test set	794 reflections (6.37%)	DCC
Wilson B-factor (Å ²)	131.3	Xtriage
Anisotropy	0.744	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 282.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	2 of 16819 reflections (0.012%)	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4399	wwPDB-VP
Average B, all atoms (Å ²)	242.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: XNM

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	0.52	3/4447 (0.1%)	0.72	4/6011 (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	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	169	TYR	N-CA	5.21	1.56	1.46
1	A	103	ASP	C-N	-5.04	1.22	1.34
1	A	168	GLY	CA-C	5.01	1.59	1.51

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	249	LEU	CA-CB-CG	6.42	130.07	115.30
1	A	228	PRO	N-CA-C	5.97	127.63	112.10
1	A	104	LEU	CB-CA-C	5.84	121.30	110.20
1	A	228	PRO	C-N-CA	5.25	134.82	121.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	232	HIS	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	4368	0	4427	318	0
2	A	31	0	22	40	0
All	All	4399	0	4449	322	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 36.

All (322) 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:A:102:GLY:HA2	2:A:676:XNM:CAT	1.25	1.64
1:A:102:GLY:HA2	2:A:676:XNM:CAS	1.27	1.60
1:A:99:CYS:O	1:A:100:GLU:CG	1.67	1.42
1:A:102:GLY:CA	2:A:676:XNM:CAT	2.12	1.26
1:A:102:GLY:CA	2:A:676:XNM:CAS	2.14	1.24
1:A:144:ARG:CD	1:A:169:TYR:O	1.89	1.20
1:A:144:ARG:HD2	1:A:169:TYR:O	1.38	1.20
1:A:235:VAL:HB	1:A:243:ILE:HA	1.20	1.15
1:A:132:ALA:CB	1:A:164:ILE:HD12	1.79	1.11
1:A:99:CYS:O	1:A:100:GLU:HG3	0.95	1.11
1:A:99:CYS:N	2:A:676:XNM:HAQ	1.67	1.10
2:A:676:XNM:HAN	2:A:676:XNM:HAZ	1.11	1.08
1:A:42:ALA:HB3	2:A:676:XNM:HAE	1.37	1.06
1:A:434:TRP:HZ3	1:A:568:ARG:HA	1.13	1.06
1:A:99:CYS:H	2:A:676:XNM:HAQ	0.96	1.05
1:A:434:TRP:CZ3	1:A:568:ARG:HA	1.91	1.04
1:A:105:ARG:CG	1:A:148:PRO:HB2	1.90	1.01
1:A:152:VAL:HG22	2:A:676:XNM:HNAK	1.23	1.00
1:A:99:CYS:H	2:A:676:XNM:CAQ	1.73	1.00
1:A:327:VAL:HG11	1:A:367:LEU:HD12	1.41	0.98
1:A:144:ARG:HB2	1:A:169:TYR:O	1.64	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:426:LEU:HB3	1:A:430:TRP:CD1	1.98	0.97
1:A:221:PRO:HA	1:A:222:PHE:HB3	1.48	0.95
1:A:102:GLY:HA2	2:A:676:XNM:HAT	1.49	0.94
1:A:42:ALA:HB3	2:A:676:XNM:CAE	1.98	0.93
1:A:42:ALA:CB	2:A:676:XNM:CAE	2.46	0.93
1:A:96:MET:CE	2:A:676:XNM:CLAA	2.53	0.93
2:A:676:XNM:CAN	2:A:676:XNM:HAZ	2.00	0.92
1:A:99:CYS:O	1:A:100:GLU:CB	2.15	0.91
1:A:105:ARG:HG2	1:A:148:PRO:HB2	1.50	0.91
1:A:132:ALA:HB1	1:A:164:ILE:HD12	1.50	0.91
1:A:98:TYR:HA	2:A:676:XNM:HAP	1.53	0.91
1:A:101:GLY:HA3	1:A:153:LEU:O	1.72	0.89
1:A:220:ARG:HE	1:A:221:PRO:HD2	1.35	0.88
1:A:73:VAL:CG2	1:A:164:ILE:HB	2.03	0.88
2:A:676:XNM:HAN	2:A:676:XNM:CAZ	2.01	0.88
1:A:144:ARG:HD2	1:A:169:TYR:C	1.93	0.88
1:A:168:GLY:O	1:A:179:CYS:SG	2.31	0.87
1:A:434:TRP:HD1	1:A:435:GLN:N	1.73	0.87
1:A:152:VAL:HG22	2:A:676:XNM:NAK	1.90	0.86
1:A:102:GLY:HA2	2:A:676:XNM:HAS	1.53	0.86
1:A:249:LEU:HB2	1:A:253:VAL:H	1.39	0.86
1:A:132:ALA:CB	1:A:164:ILE:CD1	2.55	0.85
1:A:70:HIS:HB3	1:A:135:TYR:HE1	1.41	0.84
1:A:102:GLY:CA	2:A:676:XNM:HAT	2.04	0.83
1:A:219:PHE:O	1:A:220:ARG:HB2	1.79	0.83
1:A:165:ILE:HG22	1:A:166:ASP:N	1.94	0.83
1:A:120:GLY:O	1:A:123:ARG:HB2	1.79	0.82
1:A:68:LEU:HB3	1:A:135:TYR:OH	1.80	0.82
1:A:144:ARG:CB	1:A:169:TYR:O	2.29	0.80
1:A:144:ARG:CG	1:A:169:TYR:O	2.30	0.79
1:A:222:PHE:CZ	1:A:225:ASN:HB2	2.17	0.79
1:A:73:VAL:HG22	1:A:164:ILE:HB	1.65	0.78
1:A:70:HIS:HB3	1:A:135:TYR:CE1	2.18	0.77
1:A:105:ARG:HG3	1:A:148:PRO:HB2	1.66	0.77
1:A:249:LEU:HD12	1:A:252:ALA:HA	1.64	0.77
1:A:129:ILE:HG23	1:A:164:ILE:HD11	1.67	0.77
1:A:96:MET:HE1	2:A:676:XNM:CLAA	2.22	0.77
1:A:40:GLN:HB3	1:A:98:TYR:HB3	1.68	0.76
1:A:346:ILE:HG13	1:A:347:PRO:HD2	1.66	0.76
1:A:189:LEU:HD13	1:A:190:ALA:H	1.49	0.75
1:A:260:PRO:HG3	1:A:273:LEU:HD13	1.67	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:GLY:CA	2:A:676:XNM:HAS	2.07	0.75
1:A:132:ALA:HB1	1:A:164:ILE:CD1	2.20	0.72
1:A:235:VAL:HB	1:A:243:ILE:CA	2.11	0.72
1:A:434:TRP:C	1:A:434:TRP:CD1	2.63	0.71
1:A:371:VAL:HG23	1:A:372:ILE:H	1.54	0.71
1:A:223:LEU:H	1:A:224:PRO:CD	2.04	0.71
1:A:70:HIS:HE1	1:A:132:ALA:HA	1.53	0.71
1:A:115:CYS:HB3	1:A:263:ASN:HD21	1.53	0.71
1:A:223:LEU:H	1:A:224:PRO:HD2	1.56	0.71
1:A:206:TRP:HZ2	1:A:220:ARG:HH12	1.39	0.70
1:A:187:GLN:HB3	1:A:223:LEU:HD11	1.72	0.70
1:A:165:ILE:HG22	1:A:166:ASP:H	1.55	0.69
1:A:96:MET:HE3	2:A:676:XNM:CLAA	2.30	0.69
1:A:132:ALA:HB3	1:A:164:ILE:CD1	2.22	0.68
1:A:102:GLY:O	1:A:152:VAL:HA	1.92	0.68
1:A:99:CYS:C	1:A:100:GLU:CG	2.61	0.68
1:A:571:TYR:OH	1:A:590:MET:HA	1.93	0.68
1:A:232:HIS:C	1:A:234:LYS:H	1.96	0.68
1:A:143:HIS:CD2	1:A:167:LEU:HD23	2.29	0.67
1:A:367:LEU:C	1:A:369:GLN:H	1.96	0.67
1:A:235:VAL:CB	1:A:243:ILE:HA	2.13	0.67
1:A:246:TYR:OH	1:A:259:LEU:HB3	1.94	0.67
1:A:30:LEU:O	1:A:42:ALA:HA	1.96	0.66
1:A:434:TRP:CD1	1:A:435:GLN:N	2.60	0.66
1:A:317:MET:C	1:A:319:SER:H	1.98	0.66
1:A:249:LEU:HB3	1:A:253:VAL:O	1.95	0.66
1:A:281:LEU:O	1:A:282:MET:HG2	1.94	0.66
1:A:42:ALA:HB1	2:A:676:XNM:CAE	2.24	0.66
1:A:366:PRO:HB2	1:A:368:THR:HG22	1.78	0.66
1:A:134:ARG:O	1:A:138:GLU:HG2	1.96	0.65
1:A:73:VAL:HG23	1:A:164:ILE:HB	1.79	0.65
1:A:165:ILE:CG2	1:A:166:ASP:N	2.59	0.64
1:A:165:ILE:CG2	1:A:166:ASP:H	2.10	0.64
1:A:103:ASP:HA	1:A:151:ILE:O	1.98	0.64
1:A:73:VAL:HG22	1:A:164:ILE:CB	2.27	0.64
1:A:426:LEU:O	1:A:430:TRP:HB2	1.98	0.64
1:A:323:HIS:CE1	1:A:343:ASP:HB3	2.33	0.64
1:A:205:TYR:H	1:A:205:TYR:HD2	1.47	0.63
1:A:188:TYR:O	1:A:189:LEU:HB2	1.98	0.63
1:A:117:LEU:HD22	1:A:119:GLU:CG	2.28	0.63
1:A:189:LEU:HD13	1:A:190:ALA:N	2.13	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:LEU:HD22	1:A:119:GLU:HG2	1.82	0.62
1:A:118:LYS:HA	1:A:264:HIS:O	1.99	0.62
1:A:143:HIS:HD2	1:A:167:LEU:CD2	2.12	0.62
1:A:587:SER:HB2	1:A:590:MET:HE1	1.80	0.62
1:A:221:PRO:HB3	1:A:256:SER:N	2.15	0.61
1:A:152:VAL:HG23	1:A:165:ILE:HD12	1.82	0.60
1:A:362:ASN:HA	1:A:363:SER:HB2	1.82	0.60
1:A:48:GLN:HB2	1:A:50:LEU:HG	1.83	0.60
1:A:132:ALA:HB3	1:A:164:ILE:HD12	1.75	0.60
1:A:70:HIS:CE1	1:A:72:ASN:HB2	2.37	0.59
1:A:260:PRO:HG3	1:A:273:LEU:CD1	2.32	0.59
1:A:156:GLY:H	1:A:160:LEU:HD22	1.68	0.58
1:A:102:GLY:N	2:A:676:XNM:HAS	2.18	0.58
1:A:230:GLN:O	1:A:233:GLY:N	2.34	0.58
1:A:244:VAL:HG12	1:A:278:GLN:HE21	1.68	0.58
1:A:172:GLU:O	1:A:173:LEU:HB3	2.03	0.58
1:A:148:PRO:HD2	1:A:188:TYR:HE2	1.69	0.57
1:A:59:CYS:HA	1:A:62:ILE:HG22	1.86	0.57
1:A:205:TYR:N	1:A:205:TYR:CD2	2.71	0.57
1:A:300:PHE:HB3	1:A:301:GLN:NE2	2.20	0.57
1:A:273:LEU:C	1:A:275:ARG:H	2.08	0.57
1:A:205:TYR:N	1:A:205:TYR:HD2	2.02	0.57
1:A:415:GLN:N	1:A:415:GLN:OE1	2.38	0.57
1:A:227:GLN:HB3	1:A:228:PRO:HD3	1.87	0.57
1:A:103:ASP:OD1	2:A:676:XNM:CAM	2.54	0.56
1:A:260:PRO:HD3	1:A:273:LEU:HD22	1.86	0.56
1:A:105:ARG:HD3	1:A:149:GLU:OE1	2.05	0.56
1:A:273:LEU:HG	1:A:277:LEU:HD22	1.88	0.56
1:A:152:VAL:CG2	2:A:676:XNM:CAJ	2.84	0.56
1:A:119:GLU:HB2	1:A:122:ILE:HG13	1.87	0.56
1:A:144:ARG:HD3	1:A:169:TYR:O	1.99	0.55
1:A:416:ASP:CG	1:A:417:PRO:HD3	2.27	0.55
1:A:297:VAL:HB	1:A:300:PHE:HB2	1.87	0.55
1:A:227:GLN:HE22	1:A:230:GLN:NE2	2.05	0.55
1:A:143:HIS:CD2	1:A:167:LEU:CD2	2.88	0.55
1:A:367:LEU:C	1:A:369:GLN:N	2.60	0.55
1:A:367:LEU:O	1:A:369:GLN:N	2.40	0.54
1:A:72:ASN:O	1:A:164:ILE:N	2.35	0.54
1:A:133:LEU:HA	1:A:136:LEU:HD12	1.90	0.54
1:A:628:LYS:HD3	1:A:632:VAL:HG21	1.88	0.53
1:A:99:CYS:CA	2:A:676:XNM:HAQ	2.36	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:282:MET:HB2	1:A:286:ARG:HD2	1.91	0.53
1:A:188:TYR:O	1:A:189:LEU:CB	2.56	0.53
1:A:228:PRO:N	1:A:229:VAL:HB	2.24	0.53
1:A:446:ARG:O	1:A:447:LEU:HD23	2.08	0.53
1:A:202:THR:HG21	1:A:285:GLN:H	1.74	0.53
1:A:102:GLY:H	2:A:676:XNM:HAS	1.74	0.53
1:A:260:PRO:HD3	1:A:273:LEU:CD2	2.40	0.53
1:A:105:ARG:HG2	1:A:148:PRO:CB	2.33	0.52
1:A:172:GLU:HG3	1:A:175:GLN:H	1.74	0.52
1:A:227:GLN:NE2	1:A:230:GLN:HE21	2.07	0.52
1:A:102:GLY:HA3	2:A:676:XNM:CAT	2.31	0.52
1:A:282:MET:HG3	1:A:286:ARG:NH1	2.24	0.52
1:A:223:LEU:N	1:A:224:PRO:CD	2.69	0.52
1:A:205:TYR:HB3	1:A:280:MET:HB2	1.91	0.52
1:A:42:ALA:CB	2:A:676:XNM:HAE	2.14	0.52
1:A:113:ASN:HD21	1:A:117:LEU:HD21	1.75	0.52
1:A:131:SER:HA	1:A:134:ARG:HD3	1.91	0.51
1:A:96:MET:SD	2:A:676:XNM:CLAA	3.04	0.51
1:A:333:LEU:HD11	1:A:353:LEU:HD13	1.91	0.51
1:A:341:GLN:HG3	1:A:342:GLN:N	2.25	0.51
1:A:100:GLU:HB2	1:A:155:PRO:HD2	1.93	0.51
1:A:319:SER:OG	1:A:403:LEU:HB2	2.10	0.51
1:A:433:ILE:HD12	1:A:594:LEU:HD11	1.93	0.51
1:A:605:VAL:O	1:A:608:ILE:HG22	2.11	0.50
1:A:220:ARG:HE	1:A:221:PRO:CD	2.15	0.50
1:A:422:THR:O	1:A:585:GLY:HA2	2.12	0.50
1:A:629:VAL:HG13	1:A:633:MET:HG3	1.93	0.50
1:A:223:LEU:N	1:A:224:PRO:HD2	2.25	0.50
1:A:582:ARG:O	1:A:583:THR:CB	2.60	0.50
1:A:129:ILE:HG22	1:A:133:LEU:HD13	1.93	0.50
1:A:332:ASN:O	1:A:335:ASN:N	2.43	0.50
1:A:105:ARG:HA	1:A:108:LEU:HD12	1.92	0.50
1:A:148:PRO:CD	1:A:188:TYR:CE2	2.94	0.50
1:A:147:LYS:O	1:A:150:ASN:OD1	2.30	0.50
1:A:152:VAL:HG22	2:A:676:XNM:CAJ	2.41	0.49
1:A:27:GLY:HA3	1:A:45:GLN:O	2.13	0.49
1:A:245:VAL:HG12	1:A:246:TYR:H	1.76	0.49
1:A:156:GLY:H	1:A:160:LEU:CD2	2.24	0.49
1:A:186:LEU:O	1:A:188:TYR:N	2.45	0.49
1:A:208:PHE:O	1:A:211:LEU:HB3	2.12	0.49
1:A:185:THR:HB	1:A:187:GLN:HG3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:419:ARG:HD3	1:A:587:SER:OG	2.13	0.49
1:A:152:VAL:HG21	2:A:676:XNM:CAJ	2.42	0.49
1:A:112:GLU:HG2	1:A:434:TRP:HH2	1.76	0.49
1:A:222:PHE:CD2	1:A:224:PRO:HD2	2.48	0.49
1:A:125:LEU:HA	1:A:162:HIS:CD2	2.48	0.49
1:A:304:ASP:HA	1:A:307:LEU:HB3	1.94	0.49
1:A:227:GLN:NE2	1:A:230:GLN:NE2	2.61	0.48
1:A:152:VAL:CG2	2:A:676:XNM:NAK	2.71	0.48
1:A:117:LEU:HD22	1:A:119:GLU:HG3	1.93	0.48
1:A:361:LEU:HD21	1:A:370:TYR:CD2	2.48	0.48
1:A:329:GLU:C	1:A:331:GLU:H	2.16	0.48
1:A:58:TRP:HZ3	1:A:92:PRO:HG2	1.78	0.48
1:A:148:PRO:CD	1:A:188:TYR:HE2	2.26	0.48
1:A:590:MET:CE	1:A:591:VAL:HG23	2.44	0.48
1:A:60:LEU:O	1:A:64:ILE:HG12	2.13	0.48
1:A:120:GLY:HA2	1:A:123:ARG:HH11	1.77	0.48
1:A:232:HIS:C	1:A:234:LYS:N	2.64	0.48
1:A:572:ARG:C	1:A:574:LEU:H	2.17	0.48
1:A:410:VAL:C	1:A:412:ILE:H	2.17	0.48
1:A:375:THR:HG22	1:A:376:VAL:HG23	1.96	0.48
1:A:465:LEU:HD12	1:A:623:LEU:HD22	1.96	0.48
1:A:626:SER:O	1:A:630:LYS:N	2.43	0.48
1:A:144:ARG:NH2	1:A:177:GLU:OE1	2.47	0.47
1:A:227:GLN:O	1:A:230:GLN:N	2.46	0.47
1:A:102:GLY:C	2:A:676:XNM:CAS	2.82	0.47
1:A:434:TRP:HB2	1:A:567:ALA:HB1	1.95	0.47
1:A:451:GLN:C	1:A:453:THR:H	2.18	0.47
1:A:350:GLU:HG2	1:A:391:ASP:HB2	1.96	0.47
1:A:62:ILE:HG23	1:A:63:GLN:N	2.30	0.47
1:A:134:ARG:O	1:A:135:TYR:C	2.53	0.47
1:A:235:VAL:O	1:A:243:ILE:HG12	2.15	0.47
1:A:246:TYR:CE2	1:A:258:VAL:HB	2.49	0.47
1:A:117:LEU:HB2	1:A:215:CYS:O	2.14	0.47
1:A:52:PRO:HD2	1:A:53:LYS:H	1.80	0.47
1:A:310:LYS:HD2	1:A:371:VAL:HG21	1.96	0.47
1:A:62:ILE:HG13	1:A:94:LEU:HD13	1.95	0.47
1:A:55:ARG:HD2	1:A:91:LEU:HD11	1.97	0.46
1:A:440:LEU:O	1:A:601:PHE:CE2	2.69	0.46
1:A:565:GLU:HA	1:A:568:ARG:HB2	1.97	0.46
1:A:317:MET:C	1:A:319:SER:N	2.67	0.46
1:A:101:GLY:N	1:A:155:PRO:HD3	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:VAL:O	1:A:203:VAL:N	2.48	0.46
1:A:29:VAL:HA	1:A:43:ILE:O	2.16	0.46
1:A:29:VAL:HG11	2:A:676:XNM:CAC	2.45	0.46
1:A:609:TYR:HA	1:A:612:LEU:HB3	1.97	0.46
1:A:590:MET:HG3	1:A:590:MET:H	1.38	0.46
1:A:570:LEU:C	1:A:572:ARG:H	2.19	0.46
2:A:676:XNM:CBA	2:A:676:XNM:OBE	2.64	0.45
1:A:220:ARG:HH21	1:A:221:PRO:HG2	1.82	0.45
1:A:298:GLY:O	1:A:302:ALA:HB3	2.16	0.45
1:A:52:PRO:HA	1:A:55:ARG:HH11	1.81	0.45
1:A:206:TRP:CZ2	1:A:220:ARG:NH1	2.82	0.45
1:A:74:VAL:HG13	1:A:97:GLU:CG	2.47	0.45
1:A:61:GLU:HB2	1:A:178:LEU:HD22	1.99	0.45
1:A:390:PHE:H	1:A:390:PHE:HD1	1.65	0.45
1:A:298:GLY:C	1:A:300:PHE:H	2.20	0.45
1:A:235:VAL:HG22	1:A:247:ASP:OD2	2.17	0.45
1:A:221:PRO:CA	1:A:222:PHE:HB3	2.33	0.45
1:A:428:ARG:HH21	1:A:432:GLN:HE22	1.65	0.45
1:A:332:ASN:O	1:A:333:LEU:C	2.56	0.44
1:A:115:CYS:O	1:A:263:ASN:ND2	2.51	0.44
1:A:318:VAL:O	1:A:318:VAL:HG12	2.17	0.44
1:A:371:VAL:HG23	1:A:372:ILE:N	2.29	0.44
1:A:255:PHE:CD1	1:A:255:PHE:N	2.84	0.44
1:A:448:LEU:HA	1:A:608:ILE:HD11	1.99	0.44
1:A:152:VAL:HG23	1:A:165:ILE:CD1	2.47	0.44
1:A:148:PRO:HD2	1:A:188:TYR:CE2	2.49	0.44
1:A:71:PRO:HA	1:A:163:LYS:HE3	2.00	0.44
1:A:99:CYS:O	1:A:100:GLU:HB2	2.11	0.44
1:A:170:ALA:H	1:A:178:LEU:HD12	1.83	0.44
1:A:404:PRO:HB2	1:A:405:ALA:H	1.63	0.44
1:A:363:SER:H	1:A:365:GLN:H	1.66	0.44
1:A:74:VAL:HG12	1:A:75:SER:N	2.32	0.44
1:A:428:ARG:HE	1:A:432:GLN:HE22	1.65	0.44
1:A:230:GLN:O	1:A:232:HIS:N	2.51	0.43
1:A:603:LYS:HG3	1:A:607:LEU:HD12	1.99	0.43
1:A:271:GLY:O	1:A:272:LYS:C	2.56	0.43
1:A:163:LYS:O	1:A:165:ILE:HG13	2.18	0.43
1:A:172:GLU:CG	1:A:175:GLN:H	2.30	0.43
1:A:262:PRO:HG3	1:A:432:GLN:NE2	2.34	0.43
1:A:273:LEU:O	1:A:275:ARG:N	2.51	0.43
1:A:115:CYS:HB3	1:A:263:ASN:ND2	2.29	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:ARG:HA	1:A:108:LEU:HB2	2.01	0.43
1:A:120:GLY:N	1:A:121:PRO:HD2	2.33	0.43
1:A:353:LEU:HD23	1:A:388:PHE:HB2	1.99	0.43
1:A:249:LEU:CB	1:A:253:VAL:O	2.66	0.43
1:A:104:LEU:HD23	1:A:148:PRO:HA	2.00	0.43
1:A:416:ASP:HA	1:A:591:VAL:HG22	1.99	0.43
1:A:282:MET:HB3	1:A:283:TRP:H	1.68	0.43
1:A:329:GLU:HA	1:A:366:PRO:HB3	2.00	0.43
1:A:99:CYS:N	2:A:676:XNM:CAQ	2.51	0.42
1:A:120:GLY:O	1:A:123:ARG:CB	2.59	0.42
1:A:52:PRO:HA	1:A:55:ARG:NH1	2.34	0.42
1:A:139:ASN:O	1:A:141:ILE:HG13	2.20	0.42
1:A:620:ARG:C	1:A:622:ALA:H	2.21	0.42
1:A:234:LYS:O	1:A:236:ARG:N	2.52	0.42
1:A:41:VAL:HG11	1:A:80:PRO:HG3	2.00	0.42
1:A:319:SER:C	1:A:321:ARG:N	2.73	0.42
1:A:264:HIS:ND1	1:A:264:HIS:N	2.68	0.42
1:A:96:MET:HB2	2:A:676:XNM:HAE	2.02	0.42
1:A:316:ASN:HA	1:A:388:PHE:HD2	1.84	0.42
1:A:434:TRP:CD1	1:A:435:GLN:HG2	2.55	0.42
1:A:227:GLN:O	1:A:230:GLN:HG3	2.18	0.42
1:A:222:PHE:CZ	1:A:225:ASN:CB	2.97	0.42
1:A:346:ILE:HG13	1:A:347:PRO:CD	2.45	0.42
1:A:122:ILE:O	1:A:125:LEU:HB3	2.19	0.42
1:A:430:TRP:HE3	1:A:571:TYR:HD2	1.67	0.42
1:A:463:THR:HA	1:A:466:SER:HB3	2.01	0.42
1:A:312:LEU:O	1:A:324:THR:HA	2.20	0.42
1:A:228:PRO:HB2	1:A:229:VAL:HG23	2.02	0.41
1:A:280:MET:HG3	1:A:280:MET:O	2.20	0.41
1:A:122:ILE:HG22	1:A:126:LEU:HD11	2.01	0.41
1:A:436:THR:O	1:A:440:LEU:HG	2.20	0.41
1:A:145:ASP:O	1:A:146:LEU:HB2	2.19	0.41
1:A:120:GLY:O	1:A:123:ARG:N	2.53	0.41
1:A:414:LEU:C	1:A:417:PRO:HD2	2.40	0.41
1:A:346:ILE:HA	1:A:347:PRO:HD3	1.91	0.41
1:A:321:ARG:HB3	1:A:322:VAL:H	1.57	0.41
1:A:305:SER:C	1:A:307:LEU:H	2.23	0.41
1:A:420:PRO:O	1:A:421:LEU:C	2.59	0.41
1:A:249:LEU:HB2	1:A:253:VAL:N	2.21	0.41
1:A:314:VAL:HG22	1:A:386:LEU:HD23	2.02	0.41
1:A:70:HIS:HA	1:A:71:PRO:HD3	1.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:325:TYR:HA	1:A:326:PRO:HD3	1.69	0.41
1:A:336:LEU:HB2	1:A:367:LEU:HD11	2.03	0.41
1:A:416:ASP:OD1	1:A:416:ASP:N	2.52	0.41
1:A:606:ILE:O	1:A:609:TYR:HB2	2.21	0.41
1:A:227:GLN:HE22	1:A:230:GLN:HE21	1.64	0.41
1:A:561:ASP:C	1:A:563:LEU:H	2.24	0.41
1:A:155:PRO:HA	1:A:160:LEU:HD22	2.03	0.40
1:A:220:ARG:NE	1:A:221:PRO:HD2	2.18	0.40
1:A:60:LEU:HD11	1:A:175:GLN:HB3	2.02	0.40
1:A:178:LEU:HA	1:A:178:LEU:HD23	1.81	0.40
1:A:87:ALA:HA	1:A:88:PRO:HD3	1.97	0.40
1:A:216:ILE:HD13	1:A:273:LEU:HD12	2.04	0.40
1:A:465:LEU:HD22	1:A:630:LYS:HE2	2.02	0.40
1:A:266:SER:OG	1:A:269:LEU:N	2.49	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	527/677 (78%)	329 (62%)	146 (28%)	52 (10%)	1 13

All (52) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	100	GLU
1	A	187	GLN
1	A	189	LEU
1	A	220	ARG
1	A	229	VAL
1	A	235	VAL
1	A	261	THR

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Mol	Chain	Res	Type
1	A	272	LYS
1	A	368	THR
1	A	583	THR
1	A	18	LYS
1	A	50	LEU
1	A	78	GLU
1	A	202	THR
1	A	231	TRP
1	A	265	LEU
1	A	298	GLY
1	A	404	PRO
1	A	135	TYR
1	A	227	GLN
1	A	274	GLU
1	A	296	ASN
1	A	309	LEU
1	A	330	ASN
1	A	359	LEU
1	A	363	SER
1	A	372	ILE
1	A	588	ASN
1	A	52	PRO
1	A	146	LEU
1	A	271	GLY
1	A	374	CYS
1	A	392	ASN
1	A	537	MET
1	A	573	ARG
1	A	605	VAL
1	A	621	LYS
1	A	87	ALA
1	A	221	PRO
1	A	223	LEU
1	A	284	HIS
1	A	360	ALA
1	A	630	LYS
1	A	71	PRO
1	A	262	PRO
1	A	347	PRO
1	A	120	GLY
1	A	346	ILE
1	A	530	VAL

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Mol	Chain	Res	Type
1	A	433	ILE
1	A	318	VAL
1	A	326	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	487/610 (80%)	427 (88%)	60 (12%)	6	33

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	MET
1	A	34	HIS
1	A	49	GLU
1	A	53	LYS
1	A	65	MET
1	A	70	HIS
1	A	118	LYS
1	A	135	TYR
1	A	153	LEU
1	A	186	LEU
1	A	188	TYR
1	A	189	LEU
1	A	193	LEU
1	A	199	TYR
1	A	202	THR
1	A	205	TYR
1	A	219	PHE
1	A	220	ARG
1	A	222	PHE
1	A	226	TRP
1	A	245	VAL
1	A	249	LEU
1	A	255	PHE

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Mol	Chain	Res	Type
1	A	261	THR
1	A	264	HIS
1	A	272	LYS
1	A	275	ARG
1	A	294	ASN
1	A	297	VAL
1	A	306	ILE
1	A	309	LEU
1	A	311	LEU
1	A	313	SER
1	A	321	ARG
1	A	324	THR
1	A	327	VAL
1	A	328	THR
1	A	340	LEU
1	A	343	ASP
1	A	355	GLN
1	A	361	LEU
1	A	362	ASN
1	A	388	PHE
1	A	390	PHE
1	A	416	ASP
1	A	419	ARG
1	A	422	THR
1	A	423	TYR
1	A	428	ARG
1	A	434	TRP
1	A	441	LYS
1	A	444	CYS
1	A	461	TYR
1	A	537	MET
1	A	588	ASN
1	A	590	MET
1	A	601	PHE
1	A	610	ASP
1	A	616	VAL
1	A	637	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	GLN

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Mol	Chain	Res	Type
1	A	162	HIS
1	A	196	GLN
1	A	227	GLN
1	A	230	GLN
1	A	263	ASN
1	A	278	GLN
1	A	284	HIS
1	A	301	GLN
1	A	330	ASN
1	A	355	GLN
1	A	432	GLN
1	A	435	GLN
1	A	588	ASN
1	A	599	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	XNM	A	676	-	34,34,34	3.27	5 (14%)	46,46,46	2.60	10 (21%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	XNM	A	676	-	-	0/19/29/29	0/4/4/4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	676	XNM	CBC-NBB	-16.73	1.08	1.47
2	A	676	XNM	CAG-CAH	-5.53	1.40	1.48
2	A	676	XNM	CAL-NAK	-2.43	1.35	1.40
2	A	676	XNM	CAC-CAB	2.02	1.41	1.38
2	A	676	XNM	CAV-NAW	2.87	1.41	1.34

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	676	XNM	NAR-CAJ-NAI	-7.95	118.20	126.67
2	A	676	XNM	CAP-CAQ-NAR	-4.37	118.92	123.90
2	A	676	XNM	OAQ-CAV-NAW	-3.20	117.33	122.42
2	A	676	XNM	CAF-CAG-CAD	2.15	121.61	117.55
2	A	676	XNM	CBC-NBB-CAY	2.74	118.28	111.27
2	A	676	XNM	CAG-CAH-NAI	2.94	120.05	116.13
2	A	676	XNM	CAU-CAV-NAW	3.42	123.32	118.76
2	A	676	XNM	CBC-NBB-CBA	4.54	122.90	111.27
2	A	676	XNM	CAH-NAI-CAJ	6.19	121.11	116.49
2	A	676	XNM	CAQ-NAR-CAJ	9.73	123.95	115.49

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 40 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	676	XNM	40	0

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 [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	541/677 (79%)	-0.30	5 (0%) 85 80	127, 239, 345, 435	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	32	TRP	3.6
1	A	101	GLY	3.5
1	A	33	ILE	2.8
1	A	102	GLY	2.2
1	A	475	GLU	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	XNM	A	676	31/31	0.40	0.82	2.25	262,310,325,550	0

6.5 Other polymers ⓘ

There are no such residues in this entry.