



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:04 PM GMT

PDB ID : 1IKX
Title : K103N Mutant HIV-1 Reverse Transcriptase in Complex with the Inhibitor PNU142721
Authors : Lindberg, J.; Unger, T.
Deposited on : 2001-05-07
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<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

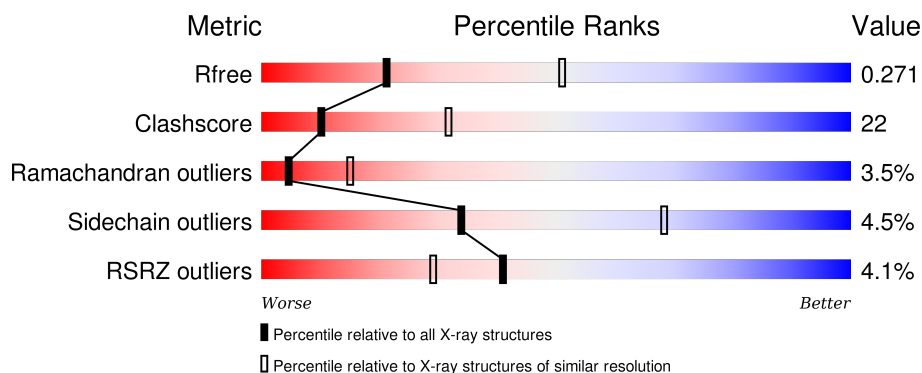
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	560	<div> <div>2%</div> <div>63%</div> <div>33%</div> <div>••</div> </div>
2	B	427	<div> <div>7%</div> <div>61%</div> <div>34%</div> <div>5%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8132 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 POL POLYPROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	557	Total	C	N	O	S	2	0	0
			4522	2925	754	835	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	103	ASN	LYS	ENGINEERED	UNP P03366
A	478	GLN	GLU	ENGINEERED	UNP P03366

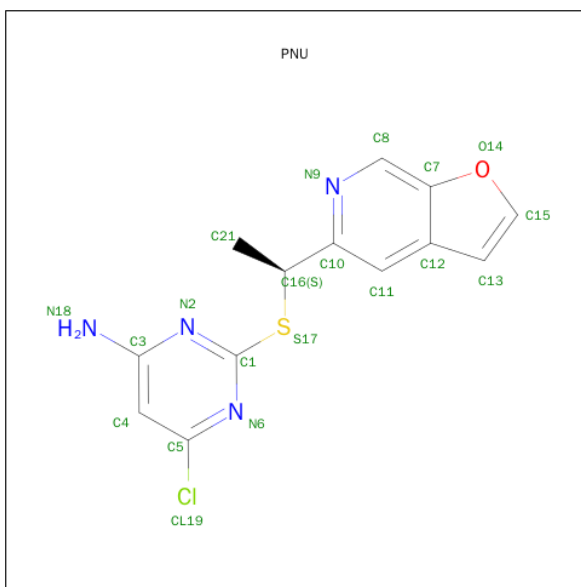
- Molecule 2 is a protein called POL POLYPROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	427	Total	C	N	O	S	33	0	0
			3529	2298	584	639	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1103	ASN	LYS	ENGINEERED	UNP P03366

- Molecule 3 is 6-CHLORO-2-(1-FURO[2,3-C]PYRIDIN-5-YL-ETHYLSULFANYL)-PYRIMIDIN-4-YLAMINE (three-letter code: PNU) (formula: C₁₃H₁₁ClN₄OS).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	S	0	0
			20	13	1	4	1	1		

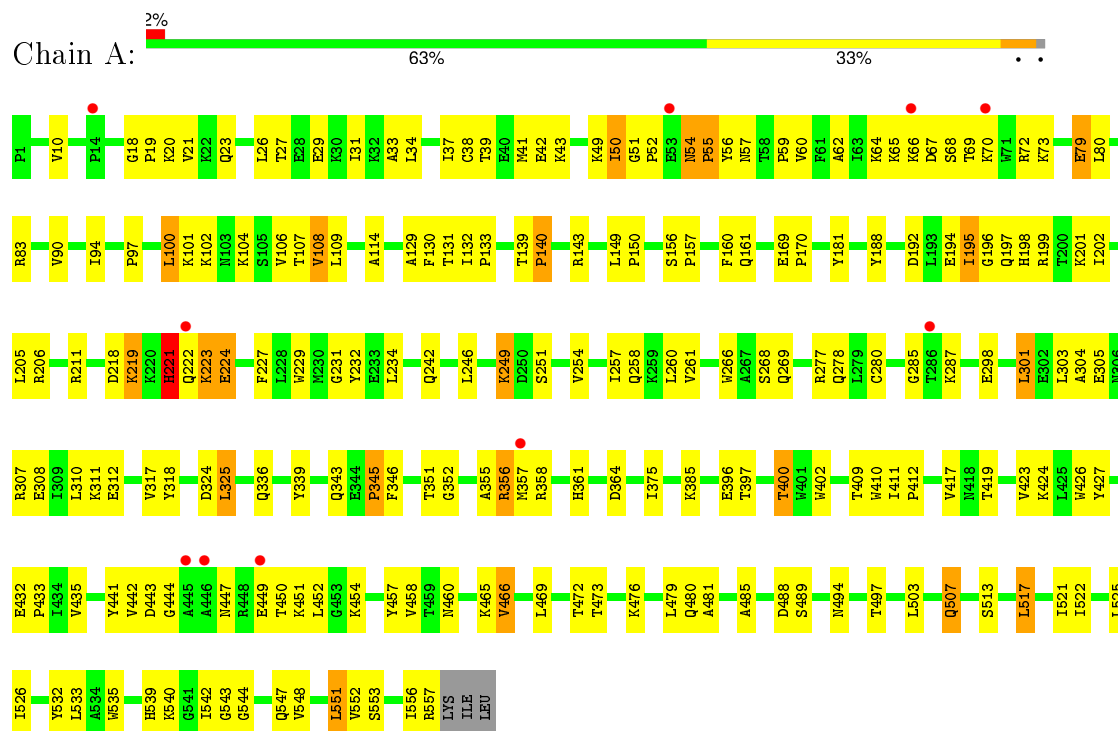
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	32	Total	O	0	0
			32	32		
4	B	29	Total	O	0	0
			29	29		

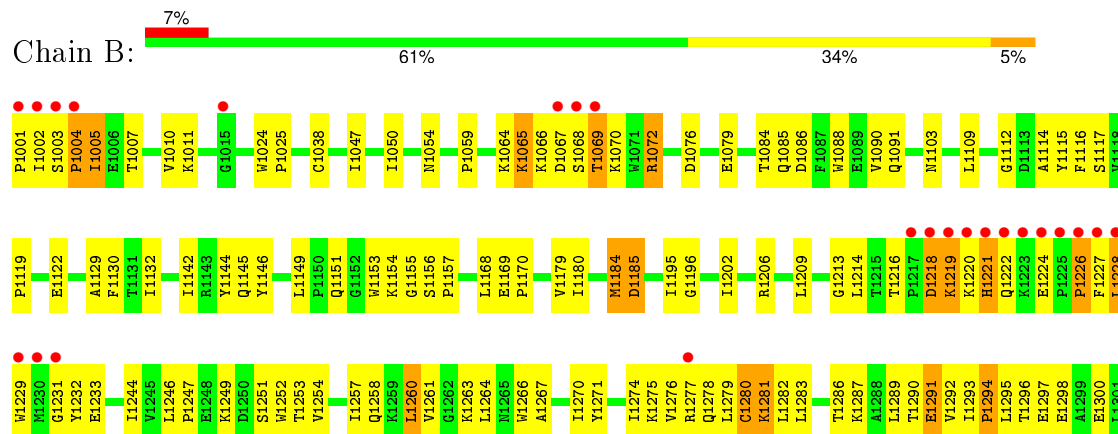
3 Residue-property plots

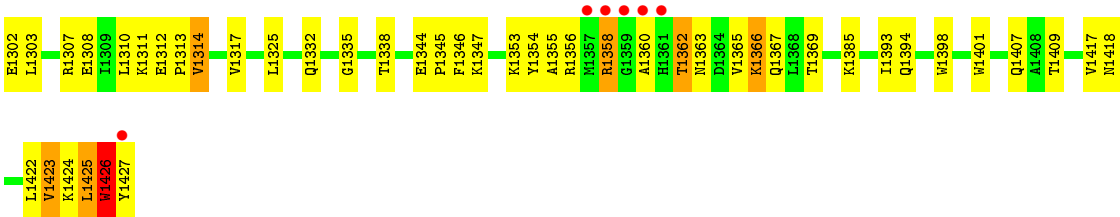
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: POL POLYPROTEIN



• Molecule 2: POL POLYPROTEIN





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	119.90Å 156.40Å 156.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	23.63 – 2.80 23.63 – 2.80	Depositor EDS
% Data completeness (in resolution range)	90.2 (23.63-2.80) 90.3 (23.63-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.69 (at 2.80Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.210 , 0.274 0.209 , 0.271	Depositor DCC
R_{free} test set	1661 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	53.5	Xtriage
Anisotropy	0.211	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 55.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 32921 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8132	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.16% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PNU

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	0.38	0/4640	0.62	0/6305
2	B	0.38	0/3634	0.63	0/4940
All	All	0.38	0/8274	0.62	0/11245

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4522	0	4572	171	0
2	B	3529	0	3559	188	0
3	A	20	0	11	0	0
4	A	32	0	0	3	0
4	B	29	0	0	3	0
All	All	8132	0	8142	352	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 22.

All (352) 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:131:THR:HG22	1:A:143:ARG:HG2	1.22	1.11
2:B:1275:LYS:HE3	2:B:1277:ARG:HB2	1.41	1.01
2:B:1066:LYS:HE2	2:B:1219:LYS:HE3	1.42	0.98
1:A:139:THR:HB	1:A:140:PRO:HD2	1.47	0.95
2:B:1422:LEU:HD23	2:B:1423:VAL:H	1.31	0.93
1:A:278:GLN:HG3	1:A:298:GLU:HB3	1.48	0.93
2:B:1266:TRP:HE1	2:B:1346:PHE:HE2	1.18	0.91
2:B:1226:PRO:HD2	2:B:1228:LEU:HD23	1.56	0.86
1:A:450:THR:HG22	1:A:452:LEU:HG	1.59	0.84
1:A:485:ALA:O	1:A:489:SER:HB2	1.79	0.82
1:A:54:ASN:HB3	1:A:143:ARG:HH21	1.44	0.82
2:B:1184:MET:O	2:B:1185:ASP:HB2	1.82	0.80
2:B:1254:VAL:HG12	2:B:1258:GLN:HE21	1.48	0.78
2:B:1365:VAL:O	2:B:1369:THR:HG23	1.84	0.78
1:A:396:GLU:O	1:A:400:THR:HG22	1.85	0.77
2:B:1279:LEU:C	2:B:1281:LYS:H	1.86	0.77
1:A:503:LEU:O	1:A:507:GLN:HB2	1.83	0.77
1:A:479:LEU:HB3	1:A:517:LEU:HD13	1.64	0.77
2:B:1007:THR:HG22	2:B:1119:PRO:HG2	1.66	0.76
2:B:1090:VAL:HG23	2:B:1091:GLN:N	2.00	0.75
1:A:450:THR:O	1:A:451:LYS:HB2	1.84	0.73
2:B:1157:PRO:HG2	2:B:1184:MET:HA	1.68	0.73
2:B:1314:VAL:O	2:B:1317:VAL:HG22	1.88	0.73
2:B:1084:THR:HG22	2:B:1088:TRP:HD1	1.52	0.73
2:B:1253:THR:O	2:B:1257:ILE:HG22	1.88	0.73
2:B:1335:GLY:HA3	2:B:1356:ARG:HG3	1.69	0.73
1:A:249:LYS:HD3	1:A:251:SER:O	1.88	0.72
2:B:1282:LEU:H	2:B:1282:LEU:HD12	1.54	0.72
2:B:1154:LYS:HE2	2:B:1184:MET:CE	2.20	0.72
1:A:356:ARG:HG2	1:A:356:ARG:HH11	1.55	0.71
1:A:268:SER:O	1:A:351:THR:HG22	1.90	0.71
2:B:1226:PRO:HG2	2:B:1227:PHE:H	1.56	0.70
1:A:21:VAL:HB	1:A:59:PRO:HD3	1.71	0.70
2:B:1047:ILE:HD12	2:B:1144:TYR:CD1	2.27	0.70
2:B:1281:LYS:HB3	2:B:1282:LEU:HD12	1.73	0.69
1:A:417:VAL:HG22	1:A:419:THR:HG23	1.74	0.69
2:B:1154:LYS:HE2	2:B:1184:MET:HE2	1.75	0.69
1:A:466:VAL:CG2	1:A:551:LEU:HD13	2.23	0.69
2:B:1224:GLU:O	2:B:1226:PRO:HD3	1.93	0.69
1:A:469:LEU:HD23	1:A:472:THR:HG21	1.74	0.69
1:A:169:GLU:HB3	1:A:170:PRO:HD3	1.75	0.69
2:B:1157:PRO:CG	2:B:1184:MET:HA	2.22	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1417:VAL:HG12	2:B:1418:ASN:N	2.08	0.68
1:A:19:PRO:HD3	1:A:80:LEU:HD13	1.76	0.68
1:A:435:VAL:HG13	2:B:1290:THR:HG21	1.76	0.68
2:B:1363:ASN:HB3	2:B:1366:LYS:HB2	1.74	0.68
1:A:131:THR:HG22	1:A:143:ARG:CG	2.14	0.68
2:B:1109:LEU:HA	2:B:1218:ASP:CG	2.13	0.67
1:A:466:VAL:HG21	1:A:551:LEU:HD13	1.74	0.67
1:A:441:TYR:O	1:A:548:VAL:HG21	1.94	0.67
2:B:1184:MET:O	2:B:1184:MET:HG3	1.95	0.67
1:A:50:ILE:HG13	1:A:51:GLY:H	1.59	0.67
1:A:131:THR:CG2	1:A:143:ARG:HH11	2.08	0.66
2:B:1059:PRO:HG2	2:B:1076:ASP:HB3	1.76	0.66
2:B:1271:TYR:O	2:B:1274:ILE:HG12	1.96	0.66
2:B:1169:GLU:HB3	2:B:1170:PRO:HD3	1.79	0.65
2:B:1090:VAL:HG23	2:B:1091:GLN:H	1.61	0.65
2:B:1066:LYS:HB2	2:B:1220:LYS:HB3	1.78	0.65
2:B:1001:PRO:HB2	2:B:1117:SER:OG	1.97	0.65
1:A:64:LYS:HD3	1:A:68:SER:OG	1.96	0.65
1:A:542:ILE:HG23	2:B:1283:LEU:HD12	1.79	0.64
2:B:1310:LEU:O	2:B:1311:LYS:HB3	1.96	0.64
1:A:107:THR:HG23	1:A:198:HIS:NE2	2.13	0.64
2:B:1246:LEU:HD12	2:B:1307:ARG:HB3	1.80	0.64
1:A:308:GLU:O	1:A:311:LYS:HG2	1.98	0.63
2:B:1358:ARG:HD2	2:B:1358:ARG:N	2.14	0.63
1:A:409:THR:HG23	4:B:3027:HOH:O	1.98	0.63
1:A:107:THR:HB	1:A:223:LYS:HB3	1.80	0.63
1:A:447:ASN:OD1	1:A:449:GLU:HG2	1.98	0.63
1:A:65:LYS:HZ1	1:A:72:ARG:HD2	1.63	0.62
1:A:37:ILE:HG22	1:A:41:MET:HE2	1.81	0.62
2:B:1065:LYS:HB3	2:B:1220:LYS:HG3	1.81	0.62
2:B:1356:ARG:NH2	2:B:1362:THR:HA	2.15	0.62
1:A:107:THR:HB	1:A:223:LYS:CB	2.30	0.62
2:B:1001:PRO:HB2	2:B:1117:SER:CB	2.30	0.61
2:B:1004:PRO:O	2:B:1005:ILE:HB	1.99	0.61
2:B:1090:VAL:CG2	2:B:1091:GLN:N	2.63	0.61
2:B:1283:LEU:HD23	2:B:1283:LEU:H	1.65	0.61
1:A:223:LYS:HG3	1:A:224:GLU:N	2.14	0.61
1:A:60:VAL:HG11	1:A:130:PHE:CD2	2.36	0.61
2:B:1085:GLN:HA	2:B:1088:TRP:NE1	2.16	0.61
1:A:50:ILE:HG13	1:A:51:GLY:N	2.15	0.61
1:A:303:LEU:O	1:A:307:ARG:HG3	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1282:LEU:HD13	2:B:1293:ILE:CG2	2.31	0.61
2:B:1417:VAL:HG12	2:B:1418:ASN:H	1.66	0.61
2:B:1072:ARG:HG3	2:B:1072:ARG:HH11	1.65	0.61
2:B:1422:LEU:O	2:B:1424:LYS:N	2.34	0.60
2:B:1090:VAL:CG2	2:B:1091:GLN:H	2.13	0.60
1:A:139:THR:HB	1:A:140:PRO:CD	2.27	0.60
2:B:1282:LEU:HD13	2:B:1293:ILE:HG21	1.82	0.60
2:B:1292:VAL:O	2:B:1293:ILE:HD12	2.00	0.60
2:B:1264:LEU:HD13	2:B:1276:VAL:HG12	1.83	0.60
2:B:1050:ILE:HD12	2:B:1054:ASN:HB3	1.83	0.60
1:A:27:THR:O	1:A:31:ILE:HG13	2.02	0.59
2:B:1282:LEU:HD22	2:B:1293:ILE:HG13	1.83	0.59
2:B:1112:GLY:C	2:B:1151:GLN:HE21	2.06	0.59
1:A:218:ASP:HB3	1:A:219:LYS:HD2	1.85	0.59
1:A:324:ASP:O	1:A:343:GLN:HG2	2.02	0.59
2:B:1422:LEU:HD23	2:B:1423:VAL:N	2.12	0.58
2:B:1279:LEU:C	2:B:1281:LYS:N	2.57	0.58
2:B:1109:LEU:HA	2:B:1218:ASP:OD1	2.03	0.58
1:A:458:VAL:HG21	1:A:547:GLN:OE1	2.04	0.58
1:A:458:VAL:HB	1:A:548:VAL:HG22	1.85	0.58
1:A:356:ARG:HG2	1:A:356:ARG:NH1	2.19	0.57
2:B:1084:THR:HG22	2:B:1088:TRP:CD1	2.36	0.57
2:B:1282:LEU:HD22	2:B:1293:ILE:HG23	1.86	0.57
2:B:1422:LEU:CD2	2:B:1423:VAL:H	2.13	0.57
2:B:1001:PRO:HD2	2:B:1116:PHE:O	2.05	0.57
1:A:37:ILE:HG22	1:A:41:MET:CE	2.35	0.57
2:B:1109:LEU:HA	2:B:1218:ASP:OD2	2.05	0.57
2:B:1409:THR:HB	4:B:3036:HOH:O	2.03	0.57
1:A:301:LEU:O	1:A:304:ALA:HB3	2.05	0.57
2:B:1219:LYS:HG2	2:B:1232:TYR:HB2	1.87	0.57
2:B:1151:GLN:O	2:B:1185:ASP:OD1	2.23	0.57
2:B:1287:LYS:HD3	2:B:1291:GLU:OE2	2.05	0.57
1:A:65:LYS:HG3	1:A:66:LYS:N	2.20	0.57
2:B:1263:LYS:HE2	2:B:1425:LEU:HA	1.87	0.56
2:B:1317:VAL:CG1	2:B:1347:LYS:HD2	2.35	0.56
2:B:1065:LYS:HD3	2:B:1221:HIS:HB2	1.86	0.56
1:A:278:GLN:CG	1:A:298:GLU:HB3	2.31	0.56
2:B:1142:ILE:H	2:B:1142:ILE:HD12	1.68	0.56
1:A:277:ARG:HB2	1:A:336:GLN:NE2	2.19	0.56
2:B:1114:ALA:HB2	2:B:1214:LEU:HD22	1.87	0.56
2:B:1356:ARG:HE	2:B:1367:GLN:HE22	1.52	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:VAL:O	1:A:258:GLN:HG3	2.06	0.56
2:B:1066:LYS:O	2:B:1067:ASP:HB2	2.06	0.56
2:B:1213:GLY:O	2:B:1214:LEU:HD23	2.05	0.56
2:B:1312:GLU:HG3	2:B:1313:PRO:HD2	1.87	0.56
2:B:1086:ASP:O	2:B:1090:VAL:HG22	2.05	0.56
2:B:1311:LYS:HG2	2:B:1311:LYS:O	2.07	0.55
1:A:20:LYS:HE2	1:A:55:PRO:HB2	1.88	0.55
1:A:194:GLU:HB2	1:A:197:GLN:OE1	2.06	0.55
2:B:1270:ILE:O	2:B:1271:TYR:HB2	2.06	0.55
1:A:521:ILE:O	1:A:525:LEU:HG	2.07	0.55
2:B:1214:LEU:O	2:B:1216:THR:HG23	2.07	0.55
2:B:1325:LEU:HD12	2:B:1385:LYS:HG3	1.89	0.55
1:A:222:GLN:O	1:A:222:GLN:HG2	2.07	0.54
1:A:317:VAL:HG23	4:A:3062:HOH:O	2.07	0.54
2:B:1282:LEU:HD12	2:B:1282:LEU:N	2.20	0.54
1:A:357:MET:HG3	1:A:358:ARG:H	1.72	0.54
2:B:1202:ILE:O	2:B:1206:ARG:HG3	2.08	0.54
1:A:497:THR:O	1:A:535:TRP:HA	2.08	0.54
1:A:65:LYS:NZ	1:A:72:ARG:HD2	2.22	0.54
1:A:131:THR:CG2	1:A:143:ARG:HG2	2.15	0.54
2:B:1064:LYS:HD2	2:B:1068:SER:O	2.08	0.54
2:B:1219:LYS:HA	2:B:1232:TYR:HB2	1.89	0.54
2:B:1154:LYS:HE2	2:B:1184:MET:HE1	1.90	0.53
1:A:107:THR:HA	1:A:223:LYS:HB2	1.90	0.53
1:A:435:VAL:HG22	2:B:1290:THR:HG23	1.88	0.53
1:A:325:LEU:HD22	1:A:385:LYS:HE3	1.90	0.53
2:B:1066:LYS:HD2	2:B:1407:GLN:HE22	1.73	0.53
1:A:469:LEU:CD2	1:A:472:THR:HG21	2.39	0.53
1:A:206:ARG:NH1	1:A:218:ASP:HA	2.24	0.53
1:A:361:HIS:CD2	1:A:513:SER:OG	2.62	0.53
1:A:458:VAL:CG2	1:A:548:VAL:HG22	2.38	0.53
1:A:211:ARG:O	1:A:211:ARG:HG2	2.09	0.53
1:A:442:VAL:HG12	1:A:457:TYR:HB3	1.91	0.52
1:A:356:ARG:NH1	1:A:358:ARG:HA	2.25	0.52
1:A:277:ARG:HD3	4:A:3019:HOH:O	2.10	0.52
1:A:361:HIS:HD2	1:A:513:SER:OG	1.92	0.52
1:A:54:ASN:HB3	1:A:143:ARG:NH2	2.20	0.52
2:B:1369:THR:HG22	2:B:1398:TRP:CH2	2.45	0.52
1:A:54:ASN:HD21	1:A:56:TYR:HB2	1.73	0.52
1:A:465:LYS:HG2	1:A:466:VAL:N	2.25	0.52
1:A:54:ASN:O	1:A:56:TYR:N	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1222:GLN:HE21	2:B:1228:LEU:HD12	1.75	0.52
1:A:460:ASN:HA	2:B:1286:THR:HG22	1.92	0.51
2:B:1213:GLY:C	2:B:1214:LEU:HD23	2.30	0.51
1:A:410:TRP:CG	1:A:411:ILE:N	2.78	0.51
2:B:1422:LEU:O	2:B:1425:LEU:HG	2.09	0.51
1:A:219:LYS:HD2	1:A:219:LYS:N	2.25	0.51
1:A:231:GLY:C	1:A:242:GLN:HG2	2.31	0.51
2:B:1263:LYS:HA	2:B:1425:LEU:HD22	1.93	0.51
2:B:1276:VAL:HA	2:B:1302:GLU:OE1	2.10	0.51
1:A:469:LEU:HD21	1:A:480:GLN:HG2	1.91	0.51
2:B:1417:VAL:CG1	2:B:1418:ASN:N	2.73	0.51
2:B:1115:TYR:HB3	2:B:1149:LEU:HB2	1.92	0.51
2:B:1065:LYS:HB2	2:B:1068:SER:HB3	1.93	0.50
1:A:221:HIS:C	1:A:223:LYS:H	2.13	0.50
2:B:1065:LYS:HE2	2:B:1221:HIS:ND1	2.27	0.50
2:B:1050:ILE:HD12	2:B:1054:ASN:CB	2.40	0.50
2:B:1260:LEU:HD11	2:B:1303:LEU:HD13	1.94	0.50
2:B:1280:CYS:O	2:B:1281:LYS:O	2.29	0.50
2:B:1417:VAL:CG1	2:B:1418:ASN:H	2.24	0.50
1:A:411:ILE:HG23	1:A:411:ILE:O	2.12	0.50
2:B:1282:LEU:HD22	2:B:1293:ILE:CG1	2.42	0.50
1:A:356:ARG:HH12	1:A:358:ARG:HA	1.76	0.50
1:A:106:VAL:HG22	1:A:227:PHE:HE2	1.77	0.49
2:B:1257:ILE:HG12	2:B:1283:LEU:HD22	1.94	0.49
1:A:23:GLN:HG2	1:A:133:PRO:HG3	1.95	0.49
1:A:412:PRO:HG3	2:B:1401:TRP:CZ2	2.47	0.49
1:A:231:GLY:O	1:A:242:GLN:HG2	2.11	0.49
1:A:34:LEU:HD21	1:A:62:ALA:HB2	1.93	0.49
1:A:285:GLY:O	1:A:287:LYS:HG3	2.12	0.49
2:B:1153:TRP:CZ2	2:B:1155:GLY:HA3	2.48	0.49
1:A:305:GLU:O	1:A:308:GLU:HB3	2.13	0.49
2:B:1010:VAL:HG12	2:B:1011:LYS:N	2.28	0.49
2:B:1257:ILE:CG1	2:B:1283:LEU:HD22	2.43	0.49
2:B:1004:PRO:O	2:B:1005:ILE:CB	2.61	0.49
1:A:50:ILE:CG1	1:A:51:GLY:H	2.21	0.48
2:B:1050:ILE:CG2	2:B:1145:GLN:HG2	2.43	0.48
2:B:1149:LEU:HB3	2:B:1156:SER:OG	2.14	0.48
1:A:156:SER:HB2	1:A:157:PRO:HD3	1.94	0.48
1:A:195:ILE:HD11	1:A:199:ARG:NH2	2.28	0.48
2:B:1346:PHE:N	2:B:1346:PHE:CD1	2.82	0.48
1:A:355:ALA:O	1:A:356:ARG:C	2.52	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:LYS:C	1:A:68:SER:H	2.17	0.48
1:A:435:VAL:HG22	2:B:1290:THR:CG2	2.44	0.48
1:A:494:ASN:HB3	2:B:1289:LEU:HD12	1.95	0.48
1:A:544:GLY:O	1:A:548:VAL:HG23	2.14	0.48
2:B:1072:ARG:NH1	2:B:1072:ARG:HG3	2.27	0.48
2:B:1264:LEU:HB3	2:B:1276:VAL:HG11	1.96	0.48
1:A:57:ASN:HA	1:A:129:ALA:O	2.14	0.48
2:B:1064:LYS:HZ1	2:B:1069:THR:HA	1.78	0.48
2:B:1226:PRO:CD	2:B:1228:LEU:HD23	2.37	0.48
2:B:1278:GLN:HB3	2:B:1298:GLU:OE2	2.14	0.48
2:B:1296:THR:CG2	2:B:1298:GLU:HG2	2.44	0.48
1:A:42:GLU:OE2	1:A:49:LYS:HG3	2.13	0.47
2:B:1282:LEU:HD22	2:B:1294:PRO:HD2	1.96	0.47
2:B:1296:THR:HG22	2:B:1297:GLU:H	1.78	0.47
1:A:442:VAL:HB	1:A:481:ALA:HB1	1.97	0.47
1:A:346:PHE:N	1:A:346:PHE:CD2	2.78	0.47
2:B:1344:GLU:O	2:B:1347:LYS:HB2	2.15	0.47
2:B:1068:SER:C	2:B:1070:LYS:H	2.18	0.47
2:B:1254:VAL:O	2:B:1257:ILE:HG23	2.15	0.47
2:B:1085:GLN:HA	2:B:1088:TRP:CE2	2.48	0.47
2:B:1047:ILE:HG22	2:B:1146:TYR:HA	1.96	0.47
1:A:443:ASP:OD1	1:A:444:GLY:N	2.42	0.47
2:B:1427:TYR:O	2:B:1427:TYR:CD1	2.68	0.47
1:A:458:VAL:HG11	1:A:547:GLN:OE1	2.13	0.47
2:B:1244:ILE:HD12	2:B:1310:LEU:HD22	1.97	0.47
2:B:1195:ILE:HG23	2:B:1196:GLY:N	2.30	0.47
1:A:131:THR:CG2	1:A:143:ARG:NH1	2.76	0.47
2:B:1266:TRP:NE1	2:B:1346:PHE:HE2	2.00	0.46
1:A:79:GLU:O	1:A:83:ARG:HD2	2.15	0.46
2:B:1079:GLU:HB2	4:B:3007:HOH:O	2.15	0.46
2:B:1024:TRP:CG	2:B:1025:PRO:HD2	2.50	0.46
1:A:97:PRO:O	1:A:100:LEU:HB2	2.16	0.46
2:B:1393:ILE:HG12	2:B:1394:GLN:N	2.30	0.46
2:B:1369:THR:HG22	2:B:1398:TRP:CZ3	2.51	0.46
1:A:108:VAL:O	1:A:221:HIS:O	2.34	0.46
2:B:1335:GLY:O	2:B:1355:ALA:HA	2.16	0.46
1:A:454:LYS:O	1:A:552:VAL:HG13	2.16	0.46
1:A:479:LEU:HB3	1:A:517:LEU:CD1	2.40	0.46
2:B:1168:LEU:HD13	2:B:1180:ILE:HG21	1.97	0.46
2:B:1282:LEU:HD22	2:B:1293:ILE:CG2	2.46	0.46
1:A:198:HIS:O	1:A:202:ILE:HG12	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:GLN:HE22	1:A:60:VAL:HG22	1.81	0.46
2:B:1209:LEU:HD22	2:B:1214:LEU:HD12	1.98	0.46
1:A:38:CYS:SG	1:A:132:ILE:HD11	2.56	0.46
1:A:65:LYS:HG3	1:A:66:LYS:H	1.80	0.45
2:B:1252:TRP:CD1	2:B:1295:LEU:HD21	2.51	0.45
2:B:1279:LEU:O	2:B:1281:LYS:N	2.50	0.45
1:A:108:VAL:C	1:A:109:LEU:HD23	2.37	0.45
1:A:202:ILE:O	1:A:205:LEU:HB3	2.16	0.45
1:A:54:ASN:HD22	1:A:54:ASN:C	2.19	0.45
1:A:458:VAL:CB	1:A:548:VAL:HG22	2.46	0.45
2:B:1267:ALA:O	2:B:1270:ILE:O	2.35	0.45
1:A:90:VAL:CG1	1:A:161:GLN:HG3	2.46	0.45
1:A:361:HIS:HD2	1:A:513:SER:CB	2.30	0.45
1:A:364:ASP:HB3	1:A:423:VAL:HG13	1.98	0.45
1:A:221:HIS:C	1:A:223:LYS:N	2.70	0.45
1:A:104:LYS:HE2	1:A:194:GLU:OE1	2.16	0.45
2:B:1353:LYS:HB3	2:B:1353:LYS:NZ	2.32	0.45
2:B:1218:ASP:O	2:B:1219:LYS:CB	2.64	0.45
2:B:1066:LYS:HA	2:B:1407:GLN:NE2	2.31	0.45
1:A:139:THR:CB	1:A:140:PRO:HD2	2.28	0.45
1:A:23:GLN:CG	1:A:133:PRO:HG3	2.47	0.45
2:B:1353:LYS:HG2	2:B:1354:TYR:N	2.32	0.44
1:A:522:ILE:O	1:A:526:ILE:HG13	2.16	0.44
2:B:1332:GLN:NE2	2:B:1423:VAL:HG13	2.31	0.44
1:A:67:ASP:O	1:A:68:SER:HB3	2.17	0.44
2:B:1426:TRP:O	2:B:1427:TYR:OXT	2.35	0.44
2:B:1296:THR:O	2:B:1300:GLU:HG2	2.17	0.44
1:A:457:TYR:C	1:A:457:TYR:CD1	2.91	0.44
2:B:1153:TRP:CE2	2:B:1155:GLY:HA3	2.51	0.44
2:B:1249:LYS:HG2	2:B:1251:SER:H	1.81	0.44
2:B:1286:THR:HG22	2:B:1286:THR:O	2.16	0.44
2:B:1064:LYS:NZ	2:B:1069:THR:HA	2.32	0.44
1:A:221:HIS:O	1:A:222:GLN:HB3	2.18	0.44
1:A:317:VAL:HG22	1:A:318:TYR:N	2.32	0.43
1:A:54:ASN:ND2	1:A:56:TYR:HB2	2.33	0.43
1:A:57:ASN:OD1	1:A:131:THR:HG23	2.18	0.43
2:B:1103:ASN:HD21	2:B:1179:VAL:CG2	2.31	0.43
2:B:1103:ASN:ND2	2:B:1179:VAL:CG2	2.82	0.43
1:A:442:VAL:CG1	1:A:485:ALA:HB2	2.49	0.43
2:B:1157:PRO:HG3	2:B:1184:MET:HA	1.97	0.43
2:B:1129:ALA:HA	2:B:1144:TYR:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:THR:HG22	1:A:70:LYS:HG3	2.00	0.43
1:A:181:TYR:HB3	1:A:188:TYR:HB2	2.01	0.43
1:A:39:THR:O	1:A:43:LYS:HG3	2.19	0.43
1:A:433:PRO:HG3	1:A:532:TYR:CE2	2.54	0.43
1:A:553:SER:HA	1:A:557:ARG:CB	2.49	0.43
2:B:1332:GLN:HE22	2:B:1423:VAL:HG13	1.83	0.43
2:B:1278:GLN:HB2	2:B:1302:GLU:CG	2.49	0.43
2:B:1038:CYS:SG	2:B:1132:ILE:HD11	2.59	0.43
1:A:417:VAL:O	1:A:417:VAL:HG13	2.19	0.43
1:A:397:THR:O	1:A:400:THR:HG23	2.18	0.42
2:B:1296:THR:HG22	2:B:1297:GLU:N	2.34	0.42
1:A:195:ILE:HG12	1:A:199:ARG:CZ	2.49	0.42
1:A:364:ASP:CB	1:A:423:VAL:HG13	2.49	0.42
2:B:1425:LEU:O	2:B:1426:TRP:C	2.57	0.42
1:A:539:HIS:O	1:A:540:LYS:HD2	2.19	0.42
1:A:33:ALA:O	1:A:37:ILE:HG13	2.20	0.42
1:A:20:LYS:HG2	1:A:55:PRO:O	2.19	0.42
1:A:106:VAL:HG22	1:A:227:PHE:CE2	2.54	0.42
1:A:257:ILE:O	1:A:261:VAL:HG23	2.19	0.42
2:B:1226:PRO:HD2	2:B:1228:LEU:CD2	2.39	0.42
1:A:351:THR:CG2	1:A:352:GLY:N	2.83	0.42
2:B:1169:GLU:HB3	2:B:1170:PRO:CD	2.49	0.42
2:B:1226:PRO:CG	2:B:1227:PHE:H	2.28	0.42
1:A:339:TYR:CZ	1:A:352:GLY:HA3	2.53	0.42
2:B:1271:TYR:CG	2:B:1310:LEU:HD23	2.54	0.42
1:A:94:ILE:HG23	1:A:229:TRP:HH2	1.85	0.42
1:A:219:LYS:HD2	1:A:219:LYS:H	1.83	0.42
2:B:1122:GLU:H	2:B:1122:GLU:CD	2.23	0.42
1:A:131:THR:HG21	1:A:143:ARG:HH11	1.84	0.42
2:B:1266:TRP:CD1	2:B:1427:TYR:OH	2.68	0.42
1:A:249:LYS:HB3	4:A:3049:HOH:O	2.19	0.42
1:A:114:ALA:HB1	1:A:160:PHE:CE1	2.55	0.42
2:B:1261:VAL:HG21	2:B:1283:LEU:HD11	2.02	0.42
1:A:149:LEU:HA	1:A:150:PRO:HD3	1.90	0.42
1:A:339:TYR:CG	1:A:375:ILE:HD11	2.55	0.42
2:B:1153:TRP:CH2	2:B:1155:GLY:HA3	2.55	0.42
2:B:1066:LYS:HD2	2:B:1407:GLN:NE2	2.34	0.41
2:B:1312:GLU:HG3	2:B:1313:PRO:CD	2.50	0.41
1:A:194:GLU:O	1:A:196:GLY:N	2.53	0.41
2:B:1206:ARG:NH1	2:B:1206:ARG:HB3	2.34	0.41
2:B:1353:LYS:CG	2:B:1354:TYR:N	2.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1065:LYS:HB2	2:B:1066:LYS:H	1.74	0.41
2:B:1084:THR:CG2	2:B:1084:THR:O	2.68	0.41
2:B:1290:THR:O	2:B:1291:GLU:C	2.58	0.41
2:B:1278:GLN:HB3	2:B:1298:GLU:HB2	2.02	0.41
1:A:94:ILE:HA	1:A:94:ILE:HD13	1.73	0.41
2:B:1231:GLY:O	2:B:1232:TYR:HB3	2.21	0.41
1:A:41:MET:CE	1:A:73:LYS:HZ3	2.34	0.41
1:A:325:LEU:HD22	1:A:385:LYS:CE	2.51	0.41
1:A:426:TRP:O	1:A:427:TYR:HB3	2.20	0.41
2:B:1222:GLN:O	2:B:1222:GLN:HG3	2.19	0.41
2:B:1130:PHE:CZ	2:B:1144:TYR:HB2	2.55	0.41
1:A:65:LYS:HZ1	1:A:72:ARG:HH11	1.68	0.41
1:A:97:PRO:HG2	1:A:232:TYR:CD1	2.56	0.41
1:A:473:THR:OG1	1:A:476:LYS:HG3	2.20	0.41
2:B:1002:ILE:HG22	2:B:1003:SER:N	2.36	0.41
1:A:246:LEU:HD11	1:A:310:LEU:HD12	2.03	0.41
2:B:1219:LYS:HD3	2:B:1219:LYS:O	2.20	0.40
1:A:357:MET:HG3	1:A:358:ARG:N	2.36	0.40
2:B:1353:LYS:HB3	2:B:1353:LYS:HZ3	1.85	0.40
1:A:266:TRP:O	1:A:269:GLN:HG2	2.21	0.40
1:A:194:GLU:C	1:A:196:GLY:N	2.75	0.40
2:B:1332:GLN:HG3	2:B:1338:THR:HG23	2.03	0.40
1:A:201:LYS:HD3	1:A:201:LYS:HA	1.97	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	555/560 (99%)	491 (88%)	50 (9%)	14 (2%)	7 24
2	B	425/427 (100%)	367 (86%)	38 (9%)	20 (5%)	3 9

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	980/987 (99%)	858 (88%)	88 (9%)	34 (4%)	4	15

All (34) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	52	PRO
1	A	223	LYS
2	B	1005	ILE
2	B	1065	LYS
2	B	1184	MET
2	B	1281	LYS
2	B	1362	THR
2	B	1423	VAL
1	A	18	GLY
1	A	221	HIS
2	B	1004	PRO
2	B	1221	HIS
2	B	1426	TRP
1	A	345	PRO
1	A	356	ARG
2	B	1185	ASP
2	B	1226	PRO
2	B	1291	GLU
2	B	1360	ALA
1	A	50	ILE
1	A	55	PRO
1	A	140	PRO
2	B	1219	LYS
2	B	1280	CYS
2	B	1294	PRO
2	B	1345	PRO
2	B	1069	THR
1	A	556	ILE
2	B	1314	VAL
1	A	466	VAL
1	A	543	GLY
2	B	1247	PRO
1	A	224	GLU
1	A	195	ILE

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	495/500 (99%)	466 (94%)	29 (6%)	24	57
2	B	389/389 (100%)	378 (97%)	11 (3%)	51	84
All	All	884/889 (99%)	844 (96%)	40 (4%)	34	68

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	VAL
1	A	26	LEU
1	A	29	GLU
1	A	54	ASN
1	A	79	GLU
1	A	100	LEU
1	A	101	LYS
1	A	102	LYS
1	A	108	VAL
1	A	192	ASP
1	A	219	LYS
1	A	221	HIS
1	A	234	LEU
1	A	249	LYS
1	A	260	LEU
1	A	280	CYS
1	A	301	LEU
1	A	312	GLU
1	A	325	LEU
1	A	345	PRO
1	A	400	THR
1	A	402	TRP
1	A	424	LYS
1	A	432	GLU
1	A	488	ASP
1	A	507	GLN
1	A	517	LEU

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Mol	Chain	Res	Type
1	A	533	LEU
1	A	551	LEU
2	B	1072	ARG
2	B	1218	ASP
2	B	1228	LEU
2	B	1229	TRP
2	B	1233	GLU
2	B	1260	LEU
2	B	1308	GLU
2	B	1358	ARG
2	B	1366	LYS
2	B	1425	LEU
2	B	1426	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN
1	A	103	ASN
1	A	145	GLN
1	A	161	GLN
1	A	182	GLN
1	A	336	GLN
1	A	361	HIS
1	A	428	GLN
1	A	500	GLN
1	A	524	GLN
2	B	1103	ASN
2	B	1137	ASN
2	B	1151	GLN
2	B	1161	GLN
2	B	1208	HIS
2	B	1222	GLN
2	B	1255	ASN
2	B	1258	GLN
2	B	1332	GLN
2	B	1367	GLN
2	B	1394	GLN
2	B	1407	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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$
3	PNU	A	2000	-	16,22,22	3.29	8 (50%)	15,31,31	3.38	6 (40%)

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
3	PNU	A	2000	-	-	0/7/8/8	0/2/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2000	PNU	C5-CL19	-9.43	1.57	1.74
3	A	2000	PNU	C4-C3	2.18	1.44	1.40
3	A	2000	PNU	C4-C5	2.19	1.41	1.38
3	A	2000	PNU	C1-N2	2.49	1.38	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2000	PNU	C10-N9	2.79	1.38	1.34
3	A	2000	PNU	C3-N18	2.90	1.43	1.35
3	A	2000	PNU	C8-N9	3.37	1.37	1.32
3	A	2000	PNU	C5-N6	5.74	1.43	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2000	PNU	C4-C5-N6	-7.59	116.50	125.62
3	A	2000	PNU	N18-C3-N2	2.01	120.17	116.50
3	A	2000	PNU	C1-N2-C3	3.15	119.04	115.97
3	A	2000	PNU	C4-C5-CL19	3.68	123.64	118.84
3	A	2000	PNU	C5-N6-C1	4.67	121.48	114.62
3	A	2000	PNU	C5-C4-C3	6.91	120.32	115.39

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 [i](#)

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	557/560 (99%)	-0.29	10 (1%) 71 61	18, 49, 85, 123	2 (0%)
2	B	427/427 (100%)	-0.05	30 (7%) 19 11	19, 46, 118, 166	10 (2%)
All	All	984/987 (99%)	-0.19	40 (4%) 41 29	18, 48, 106, 166	12 (1%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	1222	GLN	11.3
2	B	1221	HIS	10.5
2	B	1226	PRO	8.4
2	B	1230	MET	7.4
2	B	1003	SER	6.9
2	B	1223	LYS	6.9
2	B	1225	PRO	6.4
2	B	1001	PRO	5.4
2	B	1359	GLY	5.3
2	B	1002	ILE	5.2
2	B	1357	MET	5.1
2	B	1229	TRP	5.1
2	B	1360	ALA	4.9
2	B	1231	GLY	4.9
2	B	1219	LYS	4.2
2	B	1224	GLU	4.0
2	B	1067	ASP	4.0
2	B	1227	PHE	3.8
2	B	1358	ARG	3.6
2	B	1427	TYR	3.5
1	A	357	MET	3.2
2	B	1004	PRO	3.2
2	B	1218	ASP	3.0
2	B	1220	LYS	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	53	GLU	3.0
1	A	445	ALA	2.9
1	A	66	LYS	2.9
2	B	1217	PRO	2.9
1	A	70	LYS	2.8
1	A	446	ALA	2.6
2	B	1361	HIS	2.4
2	B	1228	LEU	2.3
2	B	1277	ARG	2.3
1	A	449	GLU	2.1
2	B	1015	GLY	2.1
2	B	1069	THR	2.1
1	A	222	GLN	2.0
2	B	1068	SER	2.0
1	A	286	THR	2.0
1	A	14	PRO	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
3	PNU	A	2000	20/20	0.97	0.16	-0.48	24,28,32,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.