



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:27 PM UTC

PDB ID : 1BK5 / pdb_00001bk5
Title : KARYOPHERIN ALPHA FROM SACCHAROMYCES CEREVISIAE
Authors : Conti, E.; Uy, M.; Leighton, L.; Blobel, G.; Kuriyan, J.
Deposited on : 1998-07-14
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

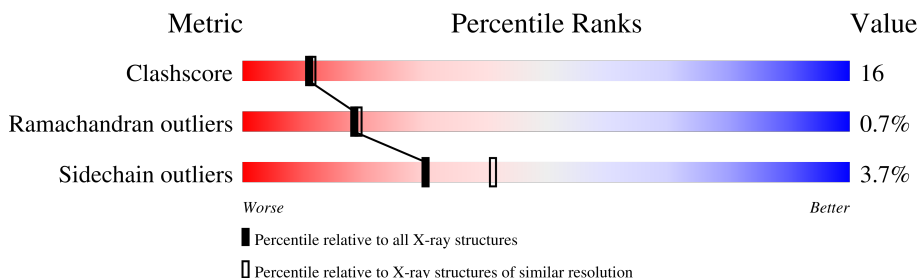
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	422	 69% 27% .
1	B	422	 64% 32% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6782 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 KARYOPHERIN ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	422	Total	C	N	O	S	88	0	0
			3275	2072	554	633	16			
1	B	422	Total	C	N	O	S	139	0	0
			3275	2072	554	633	16			

- Molecule 2 is COBALT (II) ION (CCD ID: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Co	0	0
			1	1		
2	B	1	Total	Co	0	0
			1	1		

- Molecule 3 is water.

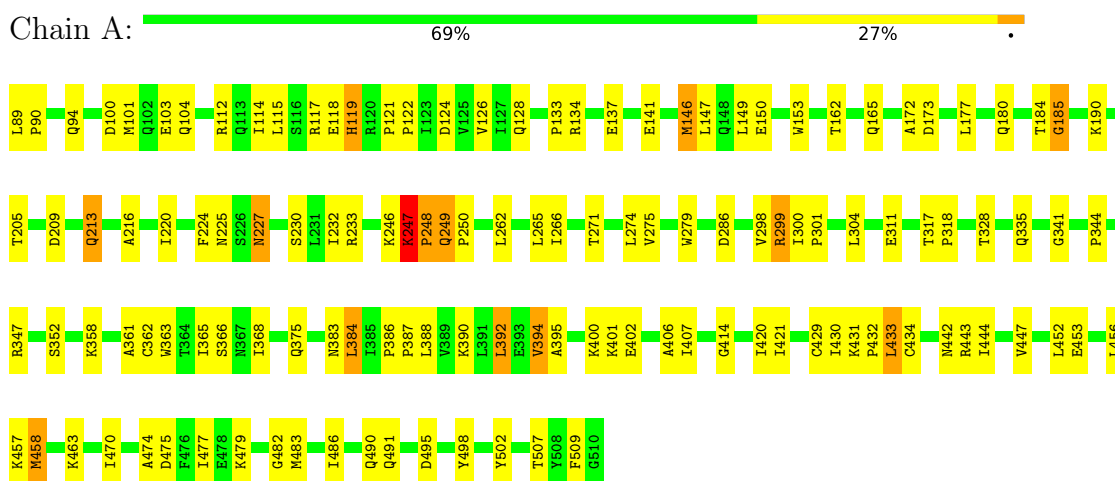
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	123	Total	O	0	0
			123	123		
3	B	107	Total	O	0	0
			107	107		

3 Residue-property plots

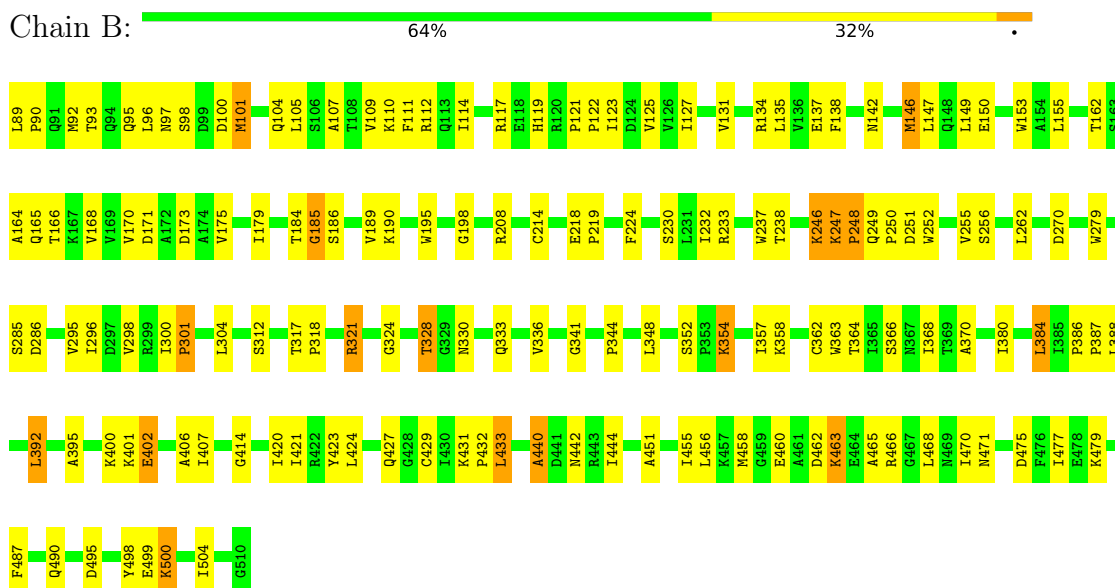
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

• Molecule 1: KARYOPHERIN ALPHA



• Molecule 1: KARYOPHERIN ALPHA



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	157.98Å 74.14Å 84.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.50 – 2.20	Depositor
% Data completeness (in resolution range)	95.0 (19.50-2.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
Refinement program	CNS 0.2	Depositor
R, R_{free}	0.237 , 0.276	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6782	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.45	0/3327	1.07	17/4526 (0.4%)
1	B	0.44	0/3327	1.05	15/4526 (0.3%)
All	All	0.45	0/6654	1.06	32/9052 (0.4%)

There are no bond length outliers.

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	247	LYS	CA-C-N	-16.73	99.65	120.23
1	A	247	LYS	C-N-CA	-16.73	99.65	120.23
1	B	247	LYS	CA-C-N	-9.95	107.40	119.84
1	B	247	LYS	C-N-CA	-9.95	107.40	119.84
1	B	247	LYS	N-CA-C	-9.30	89.27	109.81
1	B	463	LYS	N-CA-C	-8.32	101.45	111.69
1	B	368	ILE	N-CA-C	-7.75	105.62	111.90
1	A	394	VAL	N-CA-C	7.36	121.45	113.43
1	A	185	GLY	N-CA-C	7.15	120.19	112.33
1	A	328	THR	N-CA-C	-6.98	104.74	113.18
1	A	248	PRO	N-CA-C	-6.97	99.56	110.50
1	A	442	ASN	N-CA-C	6.48	119.17	111.33
1	A	184	THR	N-CA-C	6.45	119.31	111.82
1	A	507	THR	N-CA-C	6.35	118.20	111.28
1	B	186	SER	N-CA-C	-6.32	102.38	110.53
1	A	368	ILE	N-CA-C	-6.23	106.45	111.81
1	B	184	THR	N-CA-C	6.12	118.86	111.40
1	B	249	GLN	N-CA-C	-5.96	102.25	109.83
1	B	440	ALA	N-CA-C	5.93	118.36	110.53
1	B	328	THR	N-CA-C	-5.88	106.06	113.18
1	B	214	CYS	N-CA-C	-5.87	106.10	113.20
1	A	247	LYS	C-N-CD	5.72	148.44	125.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	442	ASN	N-CA-C	5.66	117.45	111.28
1	A	249	GLN	N-CA-C	-5.62	102.33	110.08
1	A	420	ILE	N-CA-C	-5.48	105.38	110.53
1	A	430	ILE	N-CA-C	5.35	115.49	110.30
1	B	142	ASN	N-CA-C	5.27	118.73	112.72
1	B	370	ALA	N-CA-C	-5.20	106.91	113.20
1	A	299	ARG	N-CA-C	5.14	118.53	111.54
1	B	312	SER	N-CA-C	5.13	117.84	109.59
1	A	141	GLU	N-CA-C	5.12	117.53	111.33
1	A	225	ASN	N-CA-C	-5.10	106.83	113.16

There are no chirality outliers.

There are no planarity outliers.

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	3275	0	3332	98	0
1	B	3275	0	3332	112	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	123	0	0	4	0
3	B	107	0	0	5	0
All	All	6782	0	6664	209	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 16.

All (209) 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:146:MET:HE1	1:A:149:LEU:HD23	1.38	1.06
1:B:101:MET:HE1	1:B:147:LEU:HD22	1.39	1.02
1:A:101:MET:HE1	1:A:147:LEU:HD22	1.43	0.96
1:A:146:MET:HA	1:A:146:MET:HE3	1.49	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:146:MET:HE3	1:B:146:MET:HA	1.47	0.92
1:B:352:SER:O	1:B:358:LYS:HE3	1.71	0.91
1:A:456:LEU:HD13	1:A:477:ILE:HD12	1.55	0.87
1:A:112:ARG:HD3	1:A:150:GLU:OE1	1.74	0.86
1:B:247:LYS:O	1:B:248:PRO:C	2.15	0.83
1:B:112:ARG:HD3	1:B:150:GLU:OE1	1.79	0.83
1:B:298:VAL:HG13	1:B:300:ILE:HG13	1.61	0.82
1:B:112:ARG:HD2	1:B:153:TRP:CE3	2.15	0.82
1:A:146:MET:CE	1:A:149:LEU:HD23	2.13	0.78
1:B:101:MET:HA	1:B:101:MET:HE3	1.67	0.76
1:A:458:MET:HA	1:A:458:MET:HE3	1.68	0.73
1:A:112:ARG:HD2	1:A:153:TRP:CE3	2.25	0.72
1:A:162:THR:OG1	1:A:165:GLN:HG3	1.93	0.69
1:A:124:ASP:O	1:A:128:GLN:HG2	1.93	0.69
1:A:352:SER:O	1:A:358:LYS:HE3	1.94	0.68
1:B:104:GLN:HE21	1:B:134:ARG:NH2	1.93	0.67
1:B:230:SER:HA	1:B:233:ARG:NH1	2.09	0.67
1:A:311:GLU:CD	1:A:311:GLU:H	2.01	0.66
1:B:246:LYS:C	1:B:247:LYS:O	2.26	0.66
1:A:347:ARG:HG3	1:A:347:ARG:HH11	1.60	0.66
1:A:224:PHE:CE2	1:A:265:LEU:HD11	2.30	0.66
1:A:133:PRO:HG2	3:A:580:HOH:O	1.95	0.65
1:B:348:LEU:HB3	3:B:564:HOH:O	1.95	0.65
1:B:122:PRO:O	1:B:125:VAL:HG12	1.95	0.65
1:B:354:LYS:HG2	1:B:357:ILE:HG13	1.78	0.65
1:B:392:LEU:HD11	1:B:433:LEU:HD13	1.80	0.64
1:B:162:THR:OG1	1:B:165:GLN:HG3	1.96	0.64
1:A:483:MET:HE3	1:A:509:PHE:HB3	1.79	0.63
1:A:341:GLY:O	1:A:344:PRO:HD2	1.99	0.62
1:B:246:LYS:HD2	1:B:286:ASP:CG	2.26	0.61
1:A:298:VAL:HG13	1:A:300:ILE:HG13	1.82	0.61
1:A:414:GLY:HA3	1:A:421:ILE:HG12	1.83	0.60
1:B:420:ILE:O	1:B:424:LEU:HD13	2.01	0.60
1:A:401:LYS:HG3	1:A:444:ILE:CD1	2.31	0.60
1:B:122:PRO:HB2	1:B:125:VAL:HG12	1.84	0.60
1:A:335:GLN:HG2	1:A:375:GLN:HE21	1.64	0.60
1:A:392:LEU:HD11	1:A:433:LEU:HD13	1.82	0.60
1:B:146:MET:HA	1:B:146:MET:CE	2.26	0.60
1:B:395:ALA:O	1:B:400:LYS:HE3	2.03	0.59
1:A:112:ARG:HD2	1:A:153:TRP:CD2	2.38	0.59
1:A:401:LYS:HG3	1:A:444:ILE:HD11	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:487:PHE:O	1:B:490:GLN:HG3	2.02	0.59
1:A:335:GLN:HG2	1:A:375:GLN:NE2	2.17	0.58
1:B:112:ARG:HD2	1:B:153:TRP:CD2	2.37	0.58
1:B:122:PRO:HB2	1:B:125:VAL:CG1	2.33	0.58
1:B:431:LYS:HB3	1:B:432:PRO:HD3	1.84	0.58
1:A:490:GLN:HE21	1:A:502:TYR:HA	1.67	0.58
1:A:347:ARG:NH1	1:A:383:ASN:O	2.37	0.57
1:A:491:GLN:HG3	3:A:549:HOH:O	2.03	0.57
1:B:495:ASP:HA	1:B:498:TYR:CE2	2.39	0.57
1:A:443:ARG:NH2	3:A:621:HOH:O	2.37	0.57
1:B:392:LEU:HD13	1:B:432:PRO:HB2	1.85	0.57
1:B:330:ASN:OD1	1:B:333:GLN:HG3	2.04	0.57
1:A:475:ASP:O	1:A:479:LYS:HG2	2.05	0.57
1:A:115:LEU:HD21	1:A:126:VAL:HG21	1.86	0.56
1:B:104:GLN:NE2	1:B:134:ARG:NH2	2.53	0.56
1:A:395:ALA:O	1:A:400:LYS:HE3	2.06	0.56
1:A:491:GLN:O	1:B:321:ARG:NH2	2.32	0.56
1:B:414:GLY:HA3	1:B:421:ILE:HG12	1.86	0.56
1:B:208:ARG:NH1	1:B:250:PRO:HD3	2.21	0.56
1:B:392:LEU:HD11	1:B:433:LEU:CD1	2.36	0.56
1:B:149:LEU:HD13	1:B:189:VAL:HA	1.87	0.56
1:B:185:GLY:O	1:B:190:LYS:HE3	2.06	0.55
1:B:392:LEU:HD12	1:B:407:ILE:HD12	1.89	0.55
1:A:114:ILE:O	1:A:117:ARG:HG2	2.07	0.55
1:B:105:LEU:O	1:B:109:VAL:HG23	2.06	0.55
1:A:100:ASP:O	1:A:104:GLN:HG3	2.07	0.55
1:A:249:GLN:NE2	3:A:534:HOH:O	2.39	0.55
1:A:300:ILE:N	1:A:301:PRO:CD	2.71	0.54
1:B:179:ILE:HG23	1:B:219:PRO:HG3	1.89	0.54
1:A:247:LYS:HA	1:A:247:LYS:CE	2.31	0.54
1:A:392:LEU:HD12	1:A:407:ILE:HD12	1.90	0.54
1:B:279:TRP:CD2	1:B:318:PRO:HB3	2.42	0.54
1:B:123:ILE:O	1:B:127:ILE:HG13	2.08	0.54
1:A:216:ALA:O	1:A:220:ILE:HG13	2.07	0.54
1:A:456:LEU:HD11	1:A:474:ALA:HA	1.90	0.54
1:B:135:LEU:HB2	1:B:155:LEU:HD21	1.88	0.54
1:A:482:GLY:O	1:A:486:ILE:HG13	2.08	0.53
1:B:300:ILE:N	1:B:301:PRO:CD	2.70	0.53
1:A:384:LEU:O	1:A:388:LEU:HB2	2.08	0.53
1:B:462:ASP:HA	1:B:465:ALA:HB3	1.90	0.53
1:B:456:LEU:HD21	1:B:477:ILE:HD12	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:247:LYS:HB3	1:B:248:PRO:HD2	1.89	0.53
1:A:101:MET:HE3	1:A:101:MET:HA	1.91	0.52
1:A:401:LYS:HE2	1:A:443:ARG:NH2	2.23	0.52
1:A:209:ASP:O	1:A:213:GLN:HB2	2.09	0.52
1:A:347:ARG:HG3	1:A:347:ARG:NH1	2.24	0.52
1:B:401:LYS:HG3	1:B:444:ILE:HG13	1.90	0.52
1:A:134:ARG:O	1:A:137:GLU:HB2	2.10	0.52
1:B:93:THR:O	1:B:97:ASN:ND2	2.43	0.52
1:A:146:MET:HA	1:A:146:MET:CE	2.30	0.51
1:A:456:LEU:HD13	1:A:477:ILE:CD1	2.34	0.51
1:A:205:THR:HG21	1:A:248:PRO:HD2	1.91	0.51
1:B:100:ASP:O	1:B:104:GLN:HG3	2.11	0.51
1:A:392:LEU:HD11	1:A:433:LEU:CD1	2.40	0.50
1:A:185:GLY:O	1:A:190:LYS:HE3	2.11	0.50
1:A:177:LEU:HA	1:A:180:GLN:HG2	1.93	0.50
1:A:172:ALA:O	1:A:173:ASP:HB2	2.10	0.50
1:B:466:ARG:HB2	1:B:468:LEU:HG	1.94	0.50
1:A:401:LYS:HE2	1:A:443:ARG:HH21	1.77	0.50
1:A:317:THR:HB	1:A:318:PRO:CD	2.41	0.50
1:A:230:SER:HA	1:A:233:ARG:HH11	1.77	0.50
1:A:443:ARG:O	1:A:447:VAL:HG23	2.13	0.49
1:B:92:MET:HE3	1:B:110:LYS:HB2	1.94	0.49
1:A:117:ARG:O	1:A:121:PRO:HD3	2.13	0.49
1:B:324:GLY:O	1:B:328:THR:HG23	2.13	0.49
1:B:101:MET:HE3	1:B:101:MET:CA	2.40	0.48
1:B:218:GLU:HB3	1:B:219:PRO:CD	2.43	0.48
1:B:500:LYS:O	1:B:504:ILE:HG13	2.13	0.48
1:A:390:LYS:O	1:A:394:VAL:HG22	2.12	0.48
1:A:463:LYS:HD2	1:A:470:ILE:N	2.29	0.48
1:B:470:ILE:HD11	1:B:475:ASP:HB2	1.93	0.48
1:A:429:CYS:O	1:A:433:LEU:HB2	2.13	0.48
1:B:208:ARG:NH2	1:B:248:PRO:O	2.34	0.48
1:B:198:GLY:HA2	1:B:238:THR:HG23	1.96	0.48
1:B:462:ASP:HA	1:B:465:ALA:CB	2.44	0.48
1:B:92:MET:HE3	1:B:110:LYS:CB	2.44	0.48
1:B:470:ILE:HG12	1:B:471:ASN:N	2.28	0.47
1:A:101:MET:HE3	1:A:101:MET:O	2.15	0.47
1:A:100:ASP:HB3	1:A:103:GLU:HB2	1.96	0.47
1:A:384:LEU:O	1:A:387:PRO:HG2	2.14	0.47
1:B:252:TRP:HA	1:B:255:VAL:HG22	1.96	0.47
1:A:453:GLU:O	1:A:457:LYS:HG3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:296:ILE:HG23	1:B:336:VAL:HG11	1.96	0.47
1:B:127:ILE:HG23	1:B:168:VAL:HG11	1.97	0.47
1:B:279:TRP:CE2	1:B:318:PRO:HB3	2.50	0.47
1:A:456:LEU:CD1	1:A:477:ILE:HD12	2.37	0.46
1:B:166:THR:O	1:B:170:VAL:HG23	2.15	0.46
1:A:90:PRO:O	1:A:94:GLN:HG3	2.16	0.46
1:A:286:ASP:OD1	1:A:286:ASP:C	2.58	0.46
1:B:362:CYS:HB2	1:B:402:GLU:HB3	1.98	0.46
1:B:498:TYR:CE1	1:B:499:GLU:HG3	2.51	0.46
1:A:347:ARG:HB2	1:A:384:LEU:HD11	1.97	0.46
1:A:452:LEU:O	1:A:456:LEU:HB2	2.15	0.46
1:B:164:ALA:HB3	3:B:546:HOH:O	2.16	0.46
1:B:384:LEU:O	1:B:388:LEU:HB2	2.15	0.46
1:B:95:GLN:O	1:B:98:SER:HB3	2.16	0.46
1:B:364:THR:HA	3:B:540:HOH:O	2.15	0.45
1:B:138:PHE:HB3	1:B:147:LEU:HG	1.97	0.45
1:B:298:VAL:O	1:B:298:VAL:HG22	2.17	0.45
1:B:386:PRO:HG3	1:B:427:GLN:OE1	2.16	0.45
1:B:363:TRP:HA	1:B:402:GLU:HG3	1.99	0.45
1:A:266:ILE:HA	1:A:274:LEU:HD12	1.98	0.45
1:B:247:LYS:O	1:B:248:PRO:O	2.34	0.45
1:A:363:TRP:HA	1:A:402:GLU:HG3	1.97	0.45
1:A:434:CYS:SG	1:A:482:GLY:HA3	2.57	0.45
1:B:121:PRO:HA	1:B:122:PRO:HD2	1.84	0.45
1:A:366:SER:HA	1:A:406:ALA:HA	1.99	0.45
1:A:279:TRP:CD2	1:A:318:PRO:HB3	2.52	0.45
1:A:483:MET:HG3	1:A:509:PHE:CD2	2.52	0.45
1:A:311:GLU:CD	1:A:311:GLU:N	2.72	0.44
1:B:101:MET:HE1	1:B:147:LEU:CD2	2.27	0.44
1:B:92:MET:HB3	1:B:111:PHE:CE1	2.53	0.44
1:B:348:LEU:CB	3:B:564:HOH:O	2.61	0.44
1:A:224:PHE:HA	1:A:232:ILE:HD12	1.99	0.44
1:B:400:LYS:NZ	3:B:582:HOH:O	2.50	0.44
1:A:431:LYS:HB3	1:A:432:PRO:HD3	1.99	0.44
1:B:440:ALA:HB1	1:B:444:ILE:HB	2.00	0.44
1:B:380:ILE:HG23	1:B:423:TYR:CE2	2.53	0.43
1:A:227:ASN:HD22	1:A:227:ASN:HA	1.57	0.43
1:A:361:ALA:O	1:A:365:ILE:HG13	2.17	0.43
1:A:247:LYS:HA	1:A:247:LYS:HD3	1.73	0.43
1:A:271:THR:O	1:A:275:VAL:HG23	2.18	0.43
1:A:421:ILE:HG21	1:A:458:MET:CB	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:GLN:CB	1:B:107:ALA:HB2	2.48	0.43
1:B:324:GLY:HA3	1:B:363:TRP:CZ3	2.54	0.43
1:A:362:CYS:HB2	1:A:402:GLU:HB3	2.00	0.43
1:A:414:GLY:HA3	1:A:421:ILE:CG1	2.49	0.43
1:B:330:ASN:CG	1:B:333:GLN:HG3	2.44	0.43
1:A:224:PHE:HA	1:A:232:ILE:CD1	2.49	0.43
1:B:460:GLU:O	1:B:463:LYS:HB3	2.19	0.43
1:A:118:GLU:O	1:A:119:HIS:O	2.37	0.42
1:B:195:TRP:CZ3	1:B:237:TRP:CZ3	3.07	0.42
1:B:101:MET:CE	1:B:147:LEU:HD22	2.28	0.42
1:B:475:ASP:O	1:B:479:LYS:HG2	2.19	0.42
1:B:134:ARG:O	1:B:137:GLU:HB2	2.20	0.42
1:B:224:PHE:HA	1:B:232:ILE:CD1	2.50	0.42
1:B:366:SER:HA	1:B:406:ALA:HA	2.01	0.42
1:B:386:PRO:HD2	1:B:387:PRO:HD2	1.99	0.42
1:B:93:THR:HG23	1:B:131:VAL:CG1	2.49	0.42
1:B:252:TRP:CE2	1:B:256:SER:HB3	2.55	0.42
1:A:386:PRO:HD2	1:A:387:PRO:HD2	2.01	0.42
1:A:121:PRO:HA	1:A:122:PRO:HD2	1.92	0.42
1:B:95:GLN:HB3	1:B:107:ALA:HB2	2.01	0.42
1:B:429:CYS:C	1:B:432:PRO:HD2	2.45	0.42
1:B:317:THR:HB	1:B:318:PRO:CD	2.51	0.41
1:B:392:LEU:HD12	1:B:407:ILE:CD1	2.50	0.41
1:B:114:ILE:O	1:B:117:ARG:HG3	2.20	0.41
1:A:363:TRP:HB2	1:A:402:GLU:HG2	2.02	0.41
1:B:96:LEU:CD2	1:B:131:VAL:HB	2.50	0.41
1:B:233:ARG:HG2	1:B:270:ASP:OD2	2.20	0.41
1:B:285:SER:OG	1:B:295:VAL:HG21	2.20	0.41
1:B:92:MET:HE2	1:B:111:PHE:CE1	2.55	0.41
1:A:249:GLN:HG3	1:A:250:PRO:HD2	2.03	0.41
1:A:458:MET:HE3	1:A:458:MET:CA	2.46	0.41
1:B:89:LEU:N	1:B:90:PRO:CD	2.84	0.41
1:B:92:MET:HE2	1:B:111:PHE:CD1	2.56	0.41
1:A:495:ASP:HA	1:A:498:TYR:CE2	2.55	0.41
1:B:451:ALA:O	1:B:455:ILE:HG13	2.20	0.41
1:A:463:LYS:CD	1:A:470:ILE:C	2.94	0.40
1:B:175:VAL:O	1:B:179:ILE:HG13	2.21	0.40
1:B:171:ASP:C	1:B:173:ASP:H	2.28	0.40
1:A:463:LYS:HG3	1:A:470:ILE:O	2.22	0.40
1:B:224:PHE:HA	1:B:232:ILE:HD12	2.03	0.40
1:B:341:GLY:O	1:B:344:PRO:HD2	2.22	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	420/422 (100%)	408 (97%)	10 (2%)	2 (0%)	24	27
1	B	420/422 (100%)	403 (96%)	13 (3%)	4 (1%)	12	11
All	All	840/844 (100%)	811 (96%)	23 (3%)	6 (1%)	18	19

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	HIS
1	B	119	HIS
1	B	248	PRO
1	A	247	LYS
1	B	185	GLY
1	B	251	ASP

5.3.2 Protein sidechains [i](#)

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	363/363 (100%)	350 (96%)	13 (4%)	31	42
1	B	363/363 (100%)	349 (96%)	14 (4%)	28	39
All	All	726/726 (100%)	699 (96%)	27 (4%)	30	41

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	89	LEU
1	A	146	MET
1	A	213	GLN
1	A	227	ASN
1	A	246	LYS
1	A	247	LYS
1	A	262	LEU
1	A	299	ARG
1	A	304	LEU
1	A	384	LEU
1	A	392	LEU
1	A	433	LEU
1	A	458	MET
1	B	101	MET
1	B	146	MET
1	B	246	LYS
1	B	262	LEU
1	B	301	PRO
1	B	304	LEU
1	B	321	ARG
1	B	354	LYS
1	B	384	LEU
1	B	392	LEU
1	B	402	GLU
1	B	433	LEU
1	B	458	MET
1	B	500	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	128	GLN
1	A	227	ASN
1	A	249	GLN
1	A	293	GLN
1	A	333	GLN
1	A	473	ASN
1	A	490	GLN
1	B	97	ASN
1	B	102	GLN
1	B	113	GLN

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Mol	Chain	Res	Type
1	B	142	ASN
1	B	192	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.