



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:56 PM UTC

PDB ID : 1EE4 / pdb_00001ee4
Title : CRYSTAL STRUCTURE OF YEAST KARYOPHERIN (IMPORTIN) ALPHA IN A COMPLEX WITH A C-MYC NLS PEPTIDE
Authors : Conti, E.
Deposited on : 2000-01-30
Resolution : 2.10 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

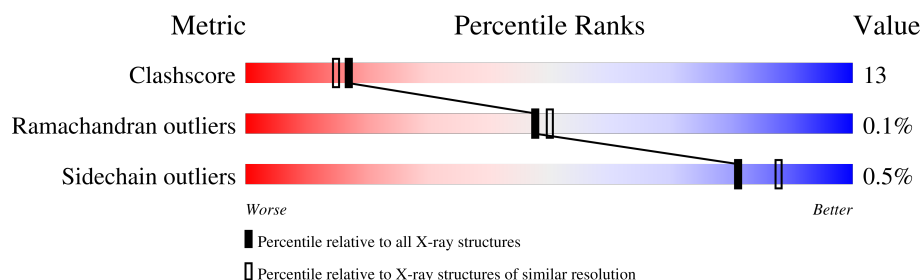
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)

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	423	
1	B	423	
2	C	9	
2	D	9	
2	E	9	
2	F	9	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6991 atoms, of which 174 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	423	Total	C	H	N	O	S	5	0	0
			3248	2012	88	528	605	15			
1	B	423	Total	C	H	N	O	S	6	0	0
			3236	2007	86	529	600	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	397	ASP	TYR	engineered mutation	UNP Q02821
B	397	ASP	TYR	engineered mutation	UNP Q02821

- Molecule 2 is a protein called MYC PROTO-ONCOGENE PROTEIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	9	Total	C	N	O	0	0	0
			69	44	14	11			
2	D	5	Total	C	N	O	0	0	0
			37	23	9	5			
2	E	9	Total	C	N	O	0	0	0
			69	44	14	11			
2	F	4	Total	C	N	O	0	0	0
			32	20	8	4			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	122	Total	O	0	0
			122	122		
3	C	14	Total	O	0	0
			14	14		
3	D	4	Total	O	0	0
			4	4		

Continued on next page...

Continued from previous page...

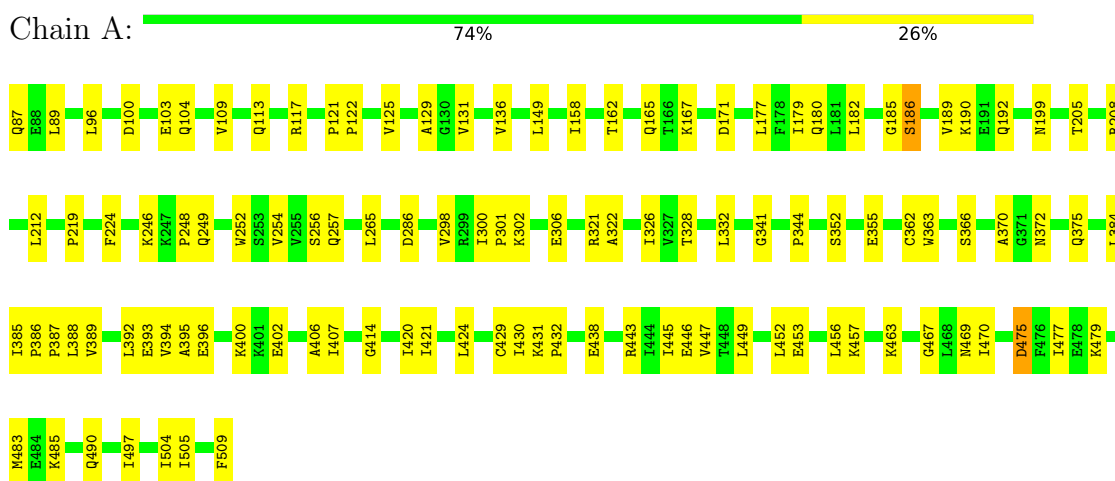
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	140	Total 140	O 140	0	0
3	E	17	Total 17	O 17	0	0
3	F	3	Total 3	O 3	0	0

3 Residue-property plots

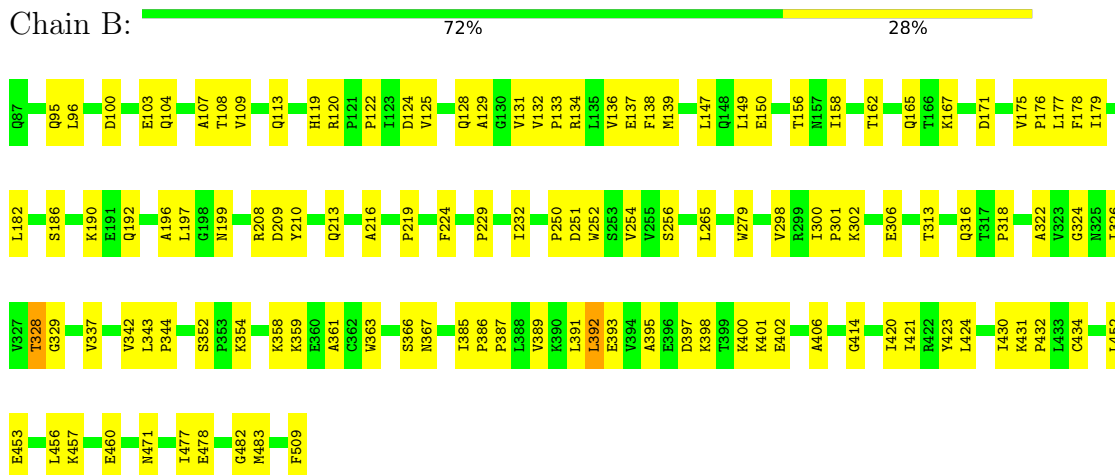
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

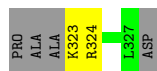
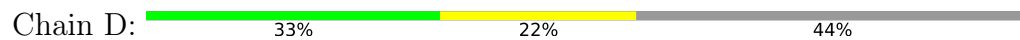


• Molecule 2: MYC PROTO-ONCOGENE PROTEIN





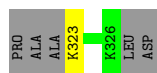
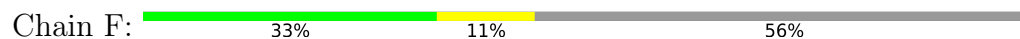
- Molecule 2: MYC PROTO-ONCOGENE PROTEIN



- Molecule 2: MYC PROTO-ONCOGENE PROTEIN



- Molecule 2: MYC PROTO-ONCOGENE PROTEIN



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	42.68Å 85.85Å 117.50Å 90.00° 93.42° 90.00°	Depositor
Resolution (Å)	30.00 – 2.10	Depositor
% Data completeness (in resolution range)	(Not available) (30.00-2.10)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.246 , 0.277	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6991	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

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	0.39	0/3211	0.92	10/4386 (0.2%)
1	B	0.39	0/3201	0.95	6/4376 (0.1%)
2	C	0.36	0/69	0.96	0/90
2	D	0.24	0/36	0.74	0/46
2	E	0.32	0/69	0.97	0/90
2	F	0.25	0/31	0.82	0/39
All	All	0.39	0/6617	0.94	16/9027 (0.2%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	328	THR	N-CA-C	-8.50	102.62	113.16
1	B	328	THR	N-CA-C	-8.11	103.39	113.28
1	B	361	ALA	N-CA-C	-7.77	102.89	111.36
1	B	430	ILE	N-CA-C	6.97	117.06	110.30
1	A	394	VAL	N-CA-C	6.06	119.02	113.20
1	A	158	ILE	N-CA-C	-6.01	104.64	110.72
1	B	158	ILE	N-CA-C	-6.00	104.46	111.00
1	B	129	ALA	N-CA-C	-5.99	105.82	113.01
1	A	430	ILE	N-CA-C	5.88	116.07	110.42
1	B	186	SER	N-CA-C	-5.59	103.15	110.53
1	A	186	SER	N-CA-C	-5.58	103.55	110.41
1	A	208	ARG	N-CA-C	-5.45	105.26	111.14
1	A	370	ALA	N-CA-C	-5.22	106.68	113.16
1	A	129	ALA	N-CA-C	-5.20	106.78	113.01
1	A	352	SER	CA-C-N	5.01	124.50	119.19
1	A	352	SER	C-N-CA	5.01	124.50	119.19

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	3160	88	3115	70	0
1	B	3150	86	3112	86	0
2	C	69	0	80	7	0
2	D	37	0	38	2	0
2	E	69	0	80	4	0
2	F	32	0	36	2	0
3	A	122	0	0	5	0
3	B	140	0	0	3	0
3	C	14	0	0	3	0
3	D	4	0	0	1	0
3	E	17	0	0	2	0
3	F	3	0	0	0	0
All	All	6817	174	6461	165	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 13.

All (165) 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:483:MET:HE2	1:B:509:PHE:HD2	1.30	0.94
1:B:483:MET:HE2	1:B:509:PHE:CD2	2.05	0.90
1:A:483:MET:HE3	1:A:509:PHE:HB3	1.59	0.85
1:A:87:GLN:HA	3:A:592:HOH:O	1.83	0.78
1:A:100:ASP:O	1:A:104:GLN:HG3	1.87	0.75
1:A:355:GLU:HG2	1:A:396:GLU:OE1	1.86	0.75
1:B:478:GLU:HB2	1:B:483:MET:HE3	1.71	0.72
1:A:122:PRO:O	1:A:125:VAL:HG12	1.88	0.72
1:B:359:LYS:HZ3	1:B:402:GLU:HG2	1.55	0.71
1:B:95:GLN:HG2	1:B:103:GLU:HB3	1.72	0.70
1:A:298:VAL:HB	3:A:600:HOH:O	1.92	0.70
1:B:359:LYS:HE2	1:B:398:LYS:O	1.92	0.69
2:E:327:LEU:O	2:E:328:ASP:HB2	1.93	0.69
1:A:117:ARG:O	1:A:121:PRO:HG3	1.94	0.67
1:B:229:PRO:HG2	3:B:573:HOH:O	1.95	0.66

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:PRO:HG2	3:A:539:HOH:O	1.96	0.66
1:A:456:LEU:HD21	1:A:477:ILE:HD12	1.79	0.65
1:A:438:GLU:OE2	1:A:485:LYS:HE3	1.96	0.65
1:A:246:LYS:HD2	1:A:286:ASP:OD2	1.98	0.64
2:C:328:ASP:C	3:C:123:HOH:O	2.40	0.64
1:B:453:GLU:HG2	1:B:457:LYS:HE3	1.77	0.64
1:A:395:ALA:O	1:A:400:LYS:HE3	1.97	0.64
1:A:205:THR:HG21	1:A:248:PRO:HD2	1.81	0.63
1:B:95:GLN:HB3	1:B:107:ALA:HB2	1.80	0.63
1:B:359:LYS:NZ	1:B:402:GLU:HG2	2.14	0.62
1:B:209:ASP:O	1:B:213:GLN:HG3	1.99	0.62
1:B:122:PRO:HB2	1:B:125:VAL:CG1	2.29	0.61
1:A:298:VAL:O	1:A:298:VAL:HG12	1.98	0.61
1:A:185:GLY:O	1:A:190:LYS:HE3	2.01	0.61
1:A:446:GLU:HG3	1:A:497:ILE:HD11	1.83	0.60
1:B:352:SER:O	1:B:358:LYS:HE3	2.01	0.60
1:A:443:ARG:O	1:A:447:VAL:HG23	2.02	0.59
1:B:431:LYS:HB3	1:B:432:PRO:HD3	1.85	0.59
1:A:149:LEU:HD12	1:A:192:GLN:HG3	1.82	0.59
1:B:298:VAL:HG13	1:B:300:ILE:HG13	1.83	0.59
1:B:100:ASP:O	1:B:104:GLN:HG3	2.01	0.59
1:B:251:ASP:OD2	1:B:254:VAL:HG23	2.03	0.59
1:B:456:LEU:HD21	1:B:477:ILE:HD12	1.85	0.58
2:C:327:LEU:O	2:C:328:ASP:HB2	2.03	0.58
1:A:300:ILE:N	1:A:301:PRO:HD2	2.18	0.58
1:A:490:GLN:NE2	1:A:505:ILE:HG13	2.20	0.57
1:B:162:THR:OG1	1:B:165:GLN:HG3	2.05	0.57
1:B:414:GLY:HA3	1:B:421:ILE:HG12	1.87	0.57
1:B:359:LYS:NZ	1:B:402:GLU:CG	2.68	0.57
1:B:395:ALA:O	1:B:400:LYS:HE3	2.05	0.56
1:A:414:GLY:HA3	1:A:421:ILE:HG12	1.88	0.56
1:A:149:LEU:HD12	1:A:192:GLN:CG	2.36	0.55
1:B:385:ILE:O	1:B:389:VAL:HG23	2.06	0.55
1:A:463:LYS:HD3	1:A:470:ILE:C	2.32	0.55
1:A:100:ASP:OD2	1:A:103:GLU:HG2	2.06	0.55
1:A:453:GLU:HG2	1:A:457:LYS:HE2	1.88	0.54
1:B:329:GLY:O	2:F:323:LYS:HE3	2.08	0.54
1:A:362:CYS:HB2	1:A:402:GLU:HB3	1.88	0.54
2:E:320:PRO:N	3:E:282:HOH:O	2.39	0.54
1:B:136:VAL:CG1	1:B:177:LEU:HD12	2.37	0.54
1:B:302:LYS:O	1:B:306:GLU:HG3	2.08	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:431:LYS:HB3	1:A:432:PRO:HD3	1.90	0.53
1:A:246:LYS:O	1:A:249:GLN:HG2	2.08	0.53
1:B:366:SER:HA	1:B:406:ALA:HA	1.90	0.53
1:A:192:GLN:HE22	2:C:326:LYS:NZ	2.07	0.53
1:B:175:VAL:HB	1:B:176:PRO:HD3	1.90	0.53
1:A:384:LEU:O	1:A:388:LEU:HB2	2.09	0.53
1:A:392:LEU:HD22	1:A:407:ILE:HD12	1.91	0.52
2:C:328:ASP:HB3	3:C:64:HOH:O	2.09	0.52
2:E:328:ASP:HB3	3:E:275:HOH:O	2.08	0.52
1:B:122:PRO:O	1:B:125:VAL:HG12	2.09	0.52
1:A:446:GLU:HG3	1:A:497:ILE:CD1	2.38	0.52
1:B:224:PHE:CE2	1:B:265:LEU:HD11	2.44	0.52
1:B:122:PRO:HB2	1:B:125:VAL:HG11	1.92	0.51
1:B:300:ILE:N	1:B:301:PRO:CD	2.73	0.51
1:A:322:ALA:O	1:A:326:ILE:HG13	2.11	0.51
1:A:212:LEU:HD12	3:A:622:HOH:O	2.11	0.51
1:B:359:LYS:HZ3	1:B:402:GLU:CG	2.25	0.50
1:B:149:LEU:HD12	1:B:192:GLN:CG	2.42	0.49
1:B:337:VAL:HG12	1:B:342:VAL:HB	1.94	0.49
1:B:397:ASP:O	1:B:401:LYS:HG3	2.13	0.49
1:A:89:LEU:CD2	1:A:125:VAL:HG22	2.43	0.48
1:B:179:ILE:HG23	1:B:219:PRO:HG3	1.95	0.48
1:B:324:GLY:O	1:B:328:THR:HG23	2.13	0.48
1:A:186:SER:OG	1:A:189:VAL:HG23	2.14	0.48
1:A:363:TRP:HA	1:A:402:GLU:HG3	1.96	0.48
1:B:208:ARG:NH1	1:B:250:PRO:HG3	2.29	0.48
1:B:359:LYS:HE2	1:B:398:LYS:C	2.38	0.48
1:A:429:CYS:O	1:A:432:PRO:HD2	2.14	0.48
1:B:478:GLU:CB	1:B:483:MET:HE3	2.43	0.48
1:B:232:ILE:HD12	3:B:635:HOH:O	2.12	0.48
1:B:322:ALA:O	1:B:326:ILE:HG13	2.14	0.48
1:A:363:TRP:CA	1:A:402:GLU:HG3	2.44	0.47
1:A:386:PRO:HD2	1:A:387:PRO:HD2	1.96	0.47
1:A:167:LYS:HG2	1:A:171:ASP:OD2	2.13	0.47
1:A:445:ILE:O	1:A:449:LEU:HG	2.14	0.47
1:B:149:LEU:HD12	1:B:192:GLN:HG2	1.96	0.47
1:B:216:ALA:O	1:B:219:PRO:HG2	2.14	0.47
1:B:343:LEU:HB2	1:B:344:PRO:HD3	1.96	0.47
1:A:96:LEU:CD2	1:A:131:VAL:HB	2.45	0.47
1:A:366:SER:HA	1:A:406:ALA:HA	1.97	0.47
1:B:359:LYS:NZ	3:B:649:HOH:O	2.48	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:279:TRP:CE2	1:B:318:PRO:HB3	2.51	0.46
1:A:254:VAL:O	1:A:257:GLN:HB2	2.15	0.46
1:A:372:ASN:OD1	1:A:375:GLN:HG3	2.15	0.46
1:A:402:GLU:OE2	2:D:324:ARG:NH2	2.45	0.46
1:A:463:LYS:HD2	1:A:469:ASN:C	2.40	0.46
1:A:89:LEU:HD21	1:A:125:VAL:HG22	1.97	0.46
1:A:179:ILE:HG23	1:A:219:PRO:HG3	1.98	0.46
1:A:224:PHE:CE2	1:A:265:LEU:HD11	2.50	0.45
1:B:167:LYS:HG2	1:B:171:ASP:OD2	2.16	0.45
1:B:108:THR:HG22	1:B:150:GLU:HB3	1.97	0.45
1:B:298:VAL:O	1:B:298:VAL:HG22	2.15	0.45
1:B:182:LEU:HD11	1:B:197:LEU:CD1	2.46	0.45
1:B:434:CYS:SG	1:B:482:GLY:HA3	2.57	0.45
1:A:341:GLY:O	1:A:344:PRO:HD2	2.18	0.44
1:A:475:ASP:O	1:A:479:LYS:HG3	2.18	0.44
3:A:616:HOH:O	2:C:322:ALA:HA	2.17	0.44
1:B:96:LEU:CD2	1:B:131:VAL:HB	2.47	0.44
1:B:313:THR:HA	1:B:316:GLN:HB2	1.99	0.44
1:B:389:VAL:O	1:B:393:GLU:HG2	2.17	0.44
1:B:386:PRO:HD2	1:B:387:PRO:HD2	1.99	0.44
1:B:452:LEU:O	1:B:456:LEU:HG	2.18	0.44
1:B:460:GLU:N	1:B:471:ASN:ND2	2.65	0.44
1:B:124:ASP:HB3	1:B:128:GLN:HE21	1.83	0.44
1:A:302:LYS:O	1:A:306:GLU:HG3	2.17	0.44
1:B:182:LEU:O	1:B:190:LYS:HE2	2.17	0.44
1:A:388:LEU:HD12	1:A:388:LEU:HA	1.87	0.43
1:A:456:LEU:CD1	1:A:504:ILE:HG23	2.48	0.43
1:B:134:ARG:O	1:B:137:GLU:HB3	2.18	0.43
1:A:392:LEU:HD23	1:A:432:PRO:HB2	1.99	0.43
1:B:175:VAL:HG21	1:B:210:TYR:CE2	2.53	0.43
1:B:132:VAL:HB	1:B:133:PRO:HD3	1.99	0.43
1:B:313:THR:HG21	1:B:354:LYS:HG3	1.99	0.43
1:A:182:LEU:O	1:A:190:LYS:HE2	2.18	0.43
1:A:199:ASN:O	2:C:323:LYS:HE2	2.18	0.43
1:B:139:MET:SD	1:B:178:PHE:CE2	3.12	0.43
1:B:136:VAL:HG12	1:B:177:LEU:HD12	2.00	0.43
1:B:385:ILE:HG13	1:B:423:TYR:HE2	1.84	0.43
1:B:119:HIS:CE1	1:B:120:ARG:HG3	2.54	0.43
1:B:175:VAL:HG21	1:B:210:TYR:HE2	1.84	0.42
1:B:367:ASN:O	2:F:323:LYS:NZ	2.52	0.42
1:B:420:ILE:O	1:B:424:LEU:HG	2.18	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:ARG:HG2	1:A:321:ARG:HH11	1.84	0.42
1:A:385:ILE:O	1:A:389:VAL:HG23	2.20	0.42
1:B:199:ASN:O	2:E:323:LYS:HE2	2.19	0.42
1:B:391:LEU:O	1:B:400:LYS:HG2	2.20	0.42
1:B:109:VAL:O	1:B:113:GLN:HG3	2.20	0.42
1:B:324:GLY:HA3	1:B:363:TRP:CZ3	2.54	0.42
1:A:177:LEU:HA	1:A:180:GLN:CG	2.49	0.42
1:B:122:PRO:HB2	1:B:125:VAL:HG12	2.02	0.42
1:B:138:PHE:HB3	1:B:147:LEU:HG	2.01	0.41
1:A:252:TRP:CE2	1:A:256:SER:HB3	2.55	0.41
1:A:136:VAL:CG1	1:A:177:LEU:HD12	2.51	0.41
2:D:323:LYS:N	3:D:119:HOH:O	2.52	0.41
1:B:385:ILE:HB	1:B:386:PRO:HD3	2.01	0.41
1:A:452:LEU:O	1:A:456:LEU:HG	2.21	0.41
1:A:420:ILE:O	1:A:424:LEU:HG	2.20	0.41
1:A:109:VAL:O	1:A:113:GLN:HG3	2.20	0.41
1:A:389:VAL:O	1:A:393:GLU:HG2	2.21	0.41
1:B:252:TRP:CE2	1:B:256:SER:HB3	2.55	0.41
1:B:279:TRP:CD2	1:B:318:PRO:HB3	2.55	0.41
1:B:156:THR:OG1	1:B:196:ALA:HB2	2.21	0.41
1:B:392:LEU:HD13	1:B:432:PRO:HB2	2.03	0.41
1:B:109:VAL:HG22	1:B:150:GLU:HG3	2.02	0.40
1:B:182:LEU:HD11	1:B:197:LEU:HD12	2.04	0.40
1:A:162:THR:OG1	1:A:165:GLN:HG3	2.21	0.40
2:C:326:LYS:NZ	3:C:94:HOH:O	2.52	0.40
1:B:104:GLN:HE21	1:B:134:ARG:NH2	2.19	0.40
1:A:298:VAL:O	1:A:298:VAL:CG1	2.68	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	421/423 (100%)	413 (98%)	7 (2%)	1 (0%)	43	44
1	B	421/423 (100%)	415 (99%)	6 (1%)	0	100	100
2	C	7/9 (78%)	7 (100%)	0	0	100	100
2	D	3/9 (33%)	3 (100%)	0	0	100	100
2	E	7/9 (78%)	7 (100%)	0	0	100	100
2	F	2/9 (22%)	2 (100%)	0	0	100	100
All	All	861/882 (98%)	847 (98%)	13 (2%)	1 (0%)	48	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	467	GLY

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	314/365 (86%)	312 (99%)	2 (1%)	78	86
1	B	326/365 (89%)	325 (100%)	1 (0%)	86	91
2	C	7/7 (100%)	7 (100%)	0	100	100
2	D	3/7 (43%)	3 (100%)	0	100	100
2	E	7/7 (100%)	7 (100%)	0	100	100
2	F	3/7 (43%)	3 (100%)	0	100	100
All	All	660/758 (87%)	657 (100%)	3 (0%)	81	88

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	332	LEU
1	A	475	ASP
1	B	392	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	104	GLN
1	A	192	GLN
1	A	213	GLN
1	A	289	GLN
1	A	490	GLN
1	B	97	ASN
1	B	104	GLN
1	B	119	HIS
1	B	128	GLN
1	B	213	GLN
1	B	215	ASN
1	B	249	GLN
1	B	257	GLN
1	B	335	GLN
1	B	339	ASN
1	B	375	GLN
1	B	488	ASN

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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

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](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.