



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 06:02 PM UTC

PDB ID : 6QNP / pdb_00006qnp
Title : CLATHRIN HEAVY CHAIN N-TERMINAL DOMAIN BOUND TO GTSE1
LIDL MOTIF
Authors : Porfetye, A.T.; Lin, Y.; Vetter, I.R.
Deposited on : 2019-02-11
Resolution : 2.70 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)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

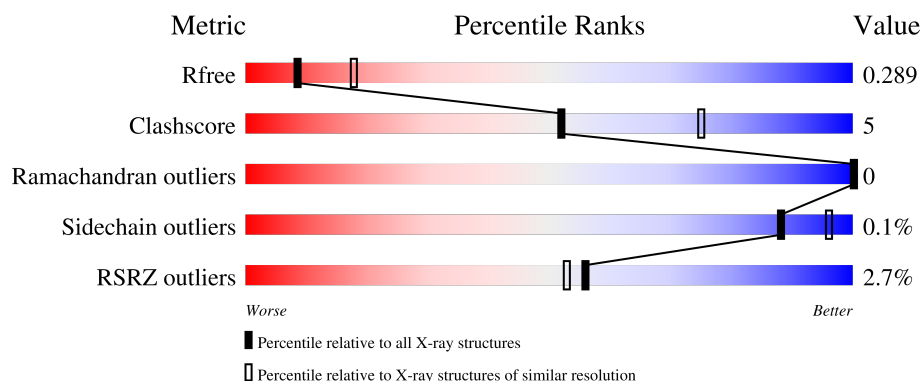
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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 $\leq 5\%$. 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	364	84% 15% .
1	B	364	91% 8% .
1	C	364	3% 84% 14% ..
1	D	364	4% 82% 15% .
2	H	67	25% 70%

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Mol	Chain	Length	Quality of chain
2	I	67	28%70%
2	J	67	12% 33%64%
2	K	67	3% 33%64%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12084 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 Clathrin heavy chain 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	361	Total	C	N	O	S	0	0	0
			2819	1792	484	525	18			
1	B	360	Total	C	N	O	S	0	0	0
			2810	1787	482	523	18			
1	C	356	Total	C	N	O	S	0	0	0
			2784	1771	479	516	18			
1	D	354	Total	C	N	O	S	0	0	0
			2770	1763	476	513	18			

- Molecule 2 is a protein called G2 and S phase-expressed protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	20	Total	C	N	O	S	0	0	0
			148	94	24	29	1			
2	I	20	Total	C	N	O	S	0	0	0
			148	94	24	29	1			
2	J	24	Total	C	N	O	S	0	0	0
			179	118	26	34	1			
2	K	24	Total	C	N	O	S	0	0	0
			180	118	26	35	1			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	86	Total	O	0	0
			86	86		
3	B	89	Total	O	0	0
			89	89		
3	C	23	Total	O	0	0
			23	23		
3	D	35	Total	O	0	0
			35	35		

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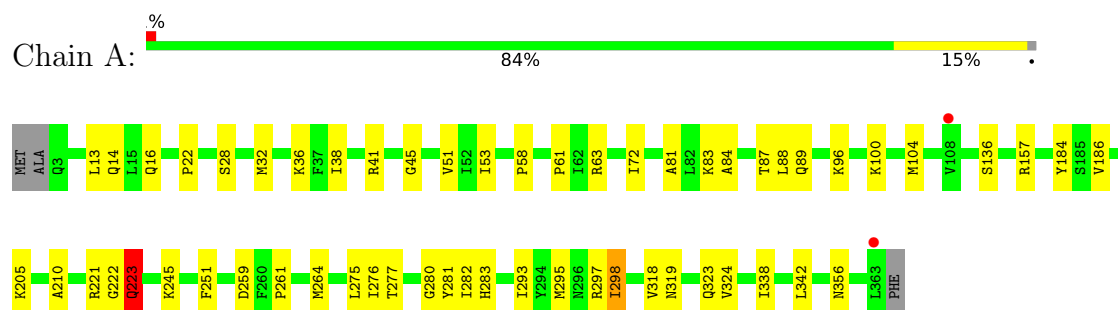
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	6	Total 6	O 6	0	0
3	I	2	Total 2	O 2	0	0
3	J	1	Total 1	O 1	0	0
3	K	4	Total 4	O 4	0	0

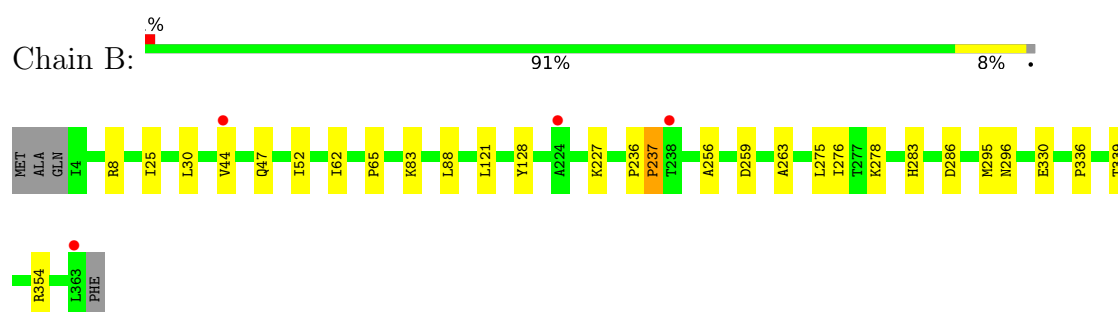
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

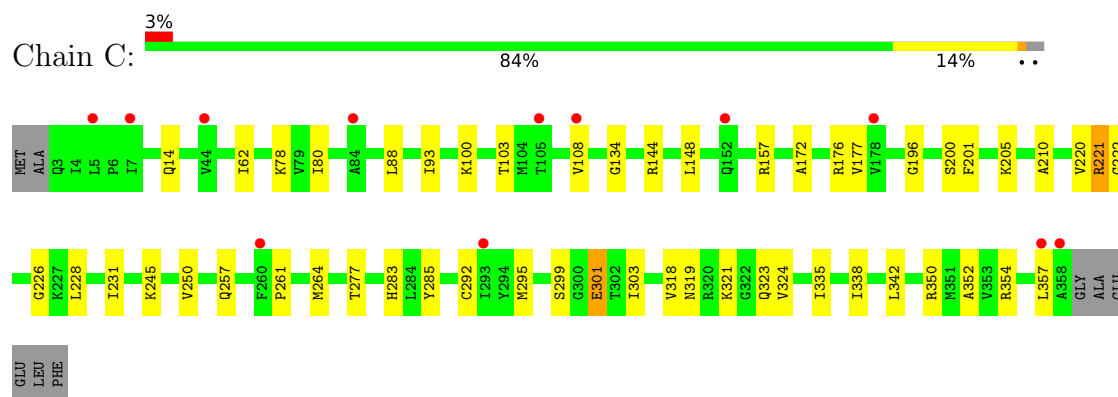
- Molecule 1: Clathrin heavy chain 1



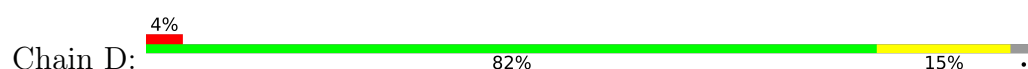
- Molecule 1: Clathrin heavy chain 1

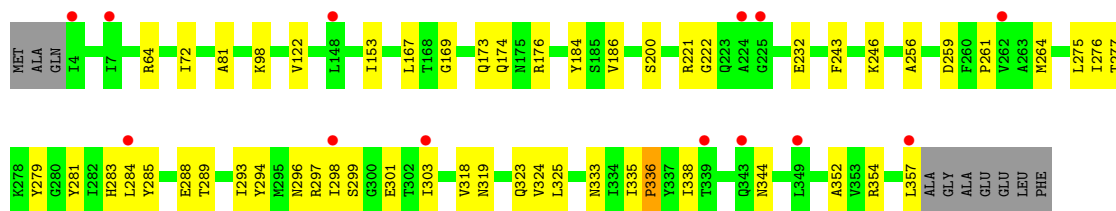


- Molecule 1: Clathrin heavy chain 1



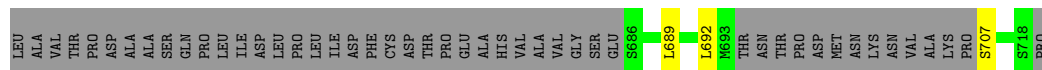
- Molecule 1: Clathrin heavy chain 1





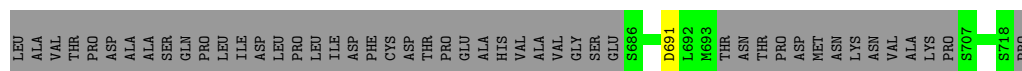
- Molecule 2: G2 and S phase-expressed protein 1

Chain H: 25% 70%



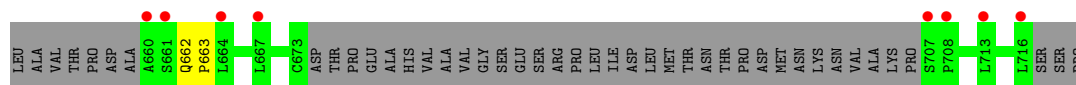
- Molecule 2: G2 and S phase-expressed protein 1

Chain I: 28% 70%



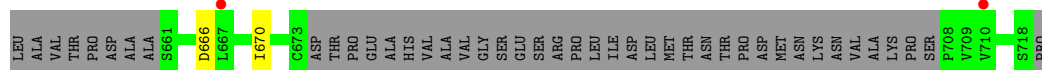
- Molecule 2: G2 and S phase-expressed protein 1

Chain J: 12% 33% 64%



- Molecule 2: G2 and S phase-expressed protein 1

Chain K: 3% 33% 64%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	95.09Å 90.15Å 100.23Å 90.00° 110.12° 90.00°	Depositor
Resolution (Å)	49.19 – 2.70 49.19 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.0 (49.19-2.70) 99.3 (49.19-2.70)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.241 , 0.291 0.245 , 0.289	Depositor DCC
R_{free} test set	2181 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	64.1	Xtriage
Anisotropy	0.416	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12084	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 74.76 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.4371e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	1/2877 (0.0%)	0.71	8/3901 (0.2%)
1	B	0.30	0/2868	0.63	2/3889 (0.1%)
1	C	0.38	0/2842	0.78	3/3854 (0.1%)
1	D	0.41	1/2828 (0.0%)	0.74	1/3835 (0.0%)
2	H	0.29	0/148	0.66	0/199
2	I	0.24	0/148	0.50	0/199
2	J	0.38	0/181	0.80	0/247
2	K	0.36	0/182	0.70	0/247
All	All	0.37	2/12074 (0.0%)	0.71	14/16371 (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
1	C	0	1
2	H	0	1
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	223	GLN	CG-CD	6.49	1.68	1.52
1	D	336	PRO	N-CD	-5.92	1.39	1.47

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	223	GLN	CG-CD-NE2	-10.08	101.28	116.40
1	C	221	ARG	CA-CB-CG	7.07	128.23	114.10
1	A	223	GLN	CB-CG-CD	7.05	124.58	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	223	GLN	CG-CD-OE1	6.43	133.66	120.80
1	A	222	GLY	CA-C-N	-6.23	111.64	122.36
1	A	222	GLY	C-N-CA	-6.23	111.64	122.36
1	D	222	GLY	N-CA-C	-6.09	98.74	113.18
1	A	223	GLN	N-CA-CB	5.96	118.51	110.59
1	A	223	GLN	CA-CB-CG	-5.66	102.78	114.10
1	C	292	CYS	CA-CB-SG	-5.36	102.08	114.40
1	B	237	PRO	CA-C-N	5.32	131.70	121.54
1	B	237	PRO	C-N-CA	5.32	131.70	121.54
1	C	221	ARG	CB-CA-C	5.18	120.31	111.46
1	A	298	ILE	CB-CA-C	-5.03	103.96	112.16

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	45	GLY	Peptide
1	C	301	GLU	Peptide
2	H	707	SER	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	2819	0	2827	38	0
1	B	2810	0	2819	16	0
1	C	2784	0	2796	35	0
1	D	2770	0	2783	32	0
2	H	148	0	157	2	0
2	I	148	0	157	1	0
2	J	179	0	185	2	0
2	K	180	0	186	2	0
3	A	86	0	0	3	0
3	B	89	0	0	2	0
3	C	23	0	0	0	0
3	D	35	0	0	0	0
3	H	6	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	I	2	0	0	0	0
3	J	1	0	0	0	0
3	K	4	0	0	0	0
All	All	12084	0	11910	122	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 5.

All (122) 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:D:173:GLN:HB2	1:D:174:GLN:OE1	1.44	1.17
1:A:223:GLN:HE21	1:A:223:GLN:HA	1.26	0.97
1:A:223:GLN:HE21	1:A:223:GLN:CA	1.77	0.96
1:A:223:GLN:HA	1:A:223:GLN:NE2	1.81	0.88
1:C:335:ILE:CG1	1:C:357:LEU:HD13	2.11	0.81
1:D:173:GLN:CB	1:D:174:GLN:OE1	2.30	0.79
1:B:44:VAL:O	1:B:47:GLN:HB2	1.89	0.72
1:D:261:PRO:HB3	1:D:275:LEU:HD11	1.78	0.65
1:C:335:ILE:HD11	1:C:357:LEU:HD13	1.78	0.64
1:B:275:LEU:HB3	1:B:283:HIS:HB2	1.80	0.63
1:D:294:TYR:HE2	1:D:296:ASN:HB2	1.62	0.63
1:D:176:ARG:NH1	1:D:221:ARG:O	2.32	0.62
1:C:14:GLN:NE2	1:C:321:LYS:O	2.34	0.60
1:B:83:LYS:HB3	1:B:88:LEU:HD23	1.83	0.60
1:D:335:ILE:HG13	1:D:336:PRO:HD3	1.82	0.60
1:C:335:ILE:CD1	1:C:357:LEU:HD13	2.30	0.60
1:C:335:ILE:HD11	1:C:357:LEU:HB2	1.85	0.57
1:C:220:VAL:HG22	1:C:222:GLY:H	1.69	0.57
1:B:286:ASP:OD1	1:B:354:ARG:NH2	2.31	0.57
1:C:338:ILE:HA	1:C:342:LEU:HD23	1.86	0.56
1:C:228:LEU:HD13	1:C:285:TYR:CD1	2.41	0.56
1:A:223:GLN:CA	1:A:223:GLN:NE2	2.51	0.55
1:A:205:LYS:NZ	1:A:210:ALA:O	2.40	0.54
1:A:223:GLN:HE21	1:A:223:GLN:C	2.14	0.54
1:C:318:VAL:HG22	1:C:324:VAL:HG22	1.90	0.54
1:C:303:ILE:HA	1:C:319:ASN:HA	1.90	0.54
1:D:284:LEU:O	1:D:293:ILE:N	2.40	0.53
1:C:144:ARG:NH2	1:C:148:LEU:O	2.41	0.52
1:C:335:ILE:HG12	1:C:357:LEU:HD13	1.89	0.52
1:A:281:TYR:HA	1:A:297:ARG:HA	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:176:ARG:NH1	1:C:222:GLY:O	2.43	0.52
1:C:261:PRO:HA	1:C:277:THR:HA	1.91	0.52
1:D:338:ILE:HD11	1:D:344:ASN:HB3	1.92	0.52
1:C:157:ARG:NH2	1:C:264:MET:O	2.38	0.51
1:D:279:TYR:HB3	1:D:297:ARG:HH21	1.75	0.51
1:B:256:ALA:HB1	1:B:259:ASP:HB2	1.92	0.50
1:A:89:GLN:HB2	2:H:689:LEU:HD12	1.94	0.50
1:A:261:PRO:HA	1:A:277:THR:HA	1.94	0.50
1:C:319:ASN:HD21	1:C:323:GLN:HB2	1.76	0.50
1:C:80:ILE:HB	1:C:93:ILE:HD11	1.94	0.50
1:A:28:SER:O	1:A:41:ARG:NH1	2.45	0.50
1:D:261:PRO:HA	1:D:277:THR:HA	1.93	0.50
1:A:251:PHE:H	1:A:295:MET:HE1	1.77	0.49
1:D:98:LYS:NZ	2:K:666:ASP:OD1	2.45	0.49
1:A:13:LEU:HD11	1:A:58:PRO:HB3	1.94	0.49
1:A:100:LYS:NZ	1:A:136:SER:O	2.41	0.49
1:C:200:SER:HB3	1:C:264:MET:HE2	1.95	0.48
1:C:88:LEU:HD11	1:C:108:VAL:HG21	1.94	0.48
1:C:103:THR:O	2:J:662:GLN:NE2	2.42	0.48
1:D:298:ILE:HD13	1:D:325:LEU:HD11	1.96	0.48
1:D:122:VAL:HG21	1:D:167:LEU:HD11	1.94	0.48
1:A:32:MET:HG3	1:A:38:ILE:HG12	1.95	0.48
1:B:65:PRO:HD2	2:I:691:ASP:HB3	1.96	0.48
1:C:205:LYS:NZ	1:C:210:ALA:O	2.47	0.48
1:C:200:SER:OG	1:C:201:PHE:N	2.47	0.47
1:D:276:ILE:HA	1:D:281:TYR:O	2.14	0.47
1:C:221:ARG:NH2	1:C:257:GLN:O	2.48	0.47
1:A:51:VAL:HG22	1:A:63:ARG:HG2	1.95	0.47
1:A:104:MET:HE3	1:A:104:MET:HB3	1.82	0.46
1:A:275:LEU:HB3	1:A:283:HIS:HB2	1.97	0.46
1:A:221:ARG:NH1	1:A:259:ASP:O	2.40	0.46
1:B:52:ILE:HB	1:B:62:ILE:HG13	1.98	0.46
1:A:83:LYS:HB2	1:A:88:LEU:HD23	1.98	0.46
1:B:263:ALA:HB3	1:B:276:ILE:HB	1.97	0.46
1:D:352:ALA:HA	1:D:357:LEU:HB2	1.98	0.46
1:A:14:GLN:NE2	3:A:421:HOH:O	2.49	0.46
1:A:72:ILE:HG13	1:A:81:ALA:HB3	1.97	0.46
1:C:352:ALA:HA	1:C:357:LEU:HD12	1.99	0.45
1:D:232:GLU:HB2	1:D:243:PHE:HB3	1.99	0.45
1:A:356:ASN:ND2	3:A:414:HOH:O	2.46	0.45
1:D:318:VAL:HG22	1:D:324:VAL:HG22	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:ILE:HA	1:A:281:TYR:O	2.17	0.45
1:B:227:LYS:NZ	3:B:417:HOH:O	2.49	0.45
1:A:319:ASN:HD21	1:A:323:GLN:HB3	1.83	0.44
1:C:172:ALA:HA	1:C:177:VAL:HA	2.00	0.44
1:A:84:ALA:O	1:A:87:THR:OG1	2.34	0.44
1:A:293:ILE:HA	1:A:342:LEU:HD11	2.00	0.44
1:B:278:LYS:NZ	3:B:412:HOH:O	2.43	0.44
1:A:36:LYS:NZ	3:A:422:HOH:O	2.50	0.44
1:A:245:LYS:HB2	1:A:245:LYS:HE2	1.80	0.44
1:B:295:MET:C	1:B:296:ASN:HD22	2.26	0.44
1:B:336:PRO:HA	1:B:339:THR:HG22	1.99	0.44
1:D:303:ILE:HG12	1:D:319:ASN:HB3	2.00	0.44
1:C:335:ILE:HG13	1:C:357:LEU:HD13	1.97	0.44
1:D:288:GLU:OE1	1:D:354:ARG:NH1	2.51	0.44
1:C:350:ARG:O	1:C:354:ARG:HB2	2.18	0.43
1:D:184:TYR:CE1	1:D:186:VAL:HG12	2.52	0.43
1:A:16:GLN:HE22	1:A:22:PRO:HD3	1.82	0.43
1:D:298:ILE:C	1:D:298:ILE:HD12	2.44	0.43
1:A:184:TYR:CE1	1:A:186:VAL:HG12	2.54	0.43
1:C:228:LEU:HD13	1:C:285:TYR:CE1	2.54	0.43
1:C:100:LYS:NZ	1:C:134:GLY:O	2.52	0.43
1:D:153:ILE:HA	1:D:169:GLY:HA2	2.01	0.43
1:D:200:SER:HB3	1:D:264:MET:HE2	1.99	0.43
1:C:299:SER:OG	1:C:301:GLU:O	2.37	0.43
1:A:83:LYS:CB	1:A:88:LEU:HD23	2.49	0.43
1:A:318:VAL:HG22	1:A:324:VAL:HG22	2.01	0.43
1:C:196:GLY:HA2	1:C:220:VAL:HB	2.01	0.43
1:D:335:ILE:HA	1:D:338:ILE:HG22	2.01	0.43
1:D:299:SER:OG	1:D:301:GLU:O	2.34	0.42
1:D:72:ILE:HG13	1:D:81:ALA:HB3	2.01	0.42
1:A:96:LYS:HA	2:H:692:LEU:HD13	2.01	0.42
1:D:246:LYS:HE2	1:D:289:THR:HA	2.00	0.42
1:A:338:ILE:HA	1:A:342:LEU:HB2	2.01	0.42
1:B:8:ARG:NH1	1:B:330:GLU:OE2	2.50	0.42
1:D:318:VAL:HA	1:D:323:GLN:O	2.19	0.42
1:A:157:ARG:NH2	1:A:264:MET:O	2.47	0.42
1:A:276:ILE:HG12	1:A:282:ILE:HD12	2.02	0.41
1:B:121:LEU:HB2	1:B:128:TYR:HB2	2.02	0.41
1:C:226:GLY:HA3	1:C:250:VAL:HB	2.01	0.41
1:A:53:ILE:HD13	1:A:61:PRO:HB3	2.02	0.41
1:A:280:GLY:O	1:A:298:ILE:N	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:283:HIS:CE1	1:C:295:MET:HE2	2.55	0.41
1:D:283:HIS:HB3	1:D:285:TYR:HE2	1.84	0.41
1:C:62:ILE:HD13	1:C:78:LYS:HZ1	1.85	0.41
1:D:333:ASN:C	1:D:336:PRO:HD2	2.46	0.41
1:B:236:PRO:HA	1:B:237:PRO:HD3	1.96	0.40
1:B:25:ILE:HG23	1:B:30:LEU:HD13	2.03	0.40
1:D:64:ARG:HB2	2:K:670:ILE:HD11	2.03	0.40
1:C:231:ILE:HG23	1:C:245:LYS:HG2	2.02	0.40
1:D:256:ALA:HB1	1:D:259:ASP:HB2	2.04	0.40
2:J:662:GLN:HA	2:J:663:PRO:HD3	1.92	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	359/364 (99%)	349 (97%)	10 (3%)	0	100	100
1	B	358/364 (98%)	342 (96%)	16 (4%)	0	100	100
1	C	354/364 (97%)	341 (96%)	13 (4%)	0	100	100
1	D	352/364 (97%)	339 (96%)	13 (4%)	0	100	100
2	H	16/67 (24%)	16 (100%)	0	0	100	100
2	I	16/67 (24%)	16 (100%)	0	0	100	100
2	J	20/67 (30%)	19 (95%)	1 (5%)	0	100	100
2	K	20/67 (30%)	19 (95%)	1 (5%)	0	100	100
All	All	1495/1724 (87%)	1441 (96%)	54 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	309/311 (99%)	308 (100%)	1 (0%)	86	94
1	B	308/311 (99%)	308 (100%)	0	100	100
1	C	306/311 (98%)	306 (100%)	0	100	100
1	D	305/311 (98%)	305 (100%)	0	100	100
2	H	19/59 (32%)	19 (100%)	0	100	100
2	I	19/59 (32%)	19 (100%)	0	100	100
2	J	22/59 (37%)	22 (100%)	0	100	100
2	K	23/59 (39%)	23 (100%)	0	100	100
All	All	1311/1480 (89%)	1310 (100%)	1 (0%)	88	96

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	223	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	60	ASN
1	A	173	GLN
1	A	223	GLN
1	A	229	HIS
1	B	47	GLN
1	B	92	ASN
1	B	102	HIS
1	B	203	GLN
1	B	296	ASN
1	B	356	ASN
1	C	3	GLN
1	C	60	ASN

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Mol	Chain	Res	Type
1	C	257	GLN
1	C	340	ASN
1	C	355	ASN
1	D	60	ASN
1	D	102	HIS
1	D	175	ASN
1	D	323	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](#)

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 [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	361/364 (99%)	0.25	2 (0%) 85 85	39, 61, 83, 109	0
1	B	360/364 (98%)	0.25	4 (1%) 78 76	39, 59, 86, 107	0
1	C	356/364 (97%)	0.62	12 (3%) 48 44	52, 83, 105, 119	0
1	D	354/364 (97%)	0.47	13 (3%) 45 41	48, 76, 107, 119	0
2	H	20/67 (29%)	0.82	0 100 100	58, 77, 103, 109	0
2	I	20/67 (29%)	0.63	0 100 100	56, 77, 98, 101	0
2	J	24/67 (35%)	1.35	8 (33%) 1 1	70, 91, 106, 107	0
2	K	24/67 (35%)	0.93	2 (8%) 17 14	65, 83, 103, 114	0
All	All	1519/1724 (88%)	0.43	41 (2%) 56 53	39, 69, 103, 119	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	7	ILE	3.7
1	C	358	ALA	3.4
1	A	363	LEU	3.1
1	B	44	VAL	3.1
1	A	108	VAL	3.0
1	D	298	ILE	2.9
2	J	664	LEU	2.9
1	B	238	THR	2.8
1	B	363	LEU	2.8
1	C	44	VAL	2.7
1	D	339	THR	2.7
1	B	224	ALA	2.6
1	D	262	VAL	2.6
1	C	152	GLN	2.6
1	D	225	GLY	2.6
2	J	667	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
2	J	660	ALA	2.5
1	C	108	VAL	2.5
1	C	260	PHE	2.5
2	J	708	PRO	2.5
1	D	349	LEU	2.4
1	D	303	ILE	2.4
2	J	707	SER	2.4
1	C	84	ALA	2.4
1	D	4	ILE	2.3
2	K	710	VAL	2.3
2	J	716	LEU	2.2
2	K	667	LEU	2.2
1	C	105	THR	2.2
1	C	5	LEU	2.2
1	D	7	ILE	2.1
1	D	357	LEU	2.1
1	D	224	ALA	2.1
2	J	661	SER	2.1
1	C	357	LEU	2.0
1	D	148	LEU	2.0
1	D	343	GLN	2.0
1	C	293	ILE	2.0
1	C	178	VAL	2.0
1	D	284	LEU	2.0
2	J	713	LEU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.