



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:40 AM UTC

PDB ID : 1DO5 / pdb_00001do5
Title : HUMAN COPPER CHAPERONE FOR SUPEROXIDE DISMUTASE DOMAIN II
Authors : Lamb, A.L.; Wernimont, A.K.; Pufahl, R.A.; O'Halloran, T.V.; Rosenzweig, A.C.
Deposited on : 1999-12-18
Resolution : 2.75 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

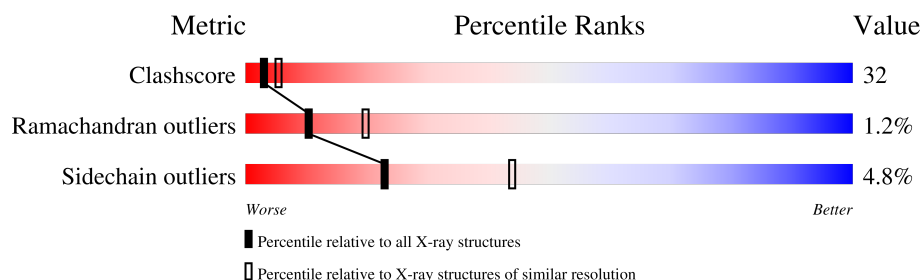
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)

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	154	
1	B	154	
1	C	154	
1	D	154	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4495 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 HUMAN COPPER CHAPERONE FOR SUPEROXIDE DISMUTASE DOMAIN II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	154	Total	C	N	O	S	0	0	0
			1143	696	218	224	5			
1	B	148	Total	C	N	O	S	0	0	0
			1094	664	210	215	5			
1	C	148	Total	C	N	O	S	0	0	0
			1094	664	210	215	5			
1	D	153	Total	C	N	O	S	0	0	0
			1134	691	216	222	5			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		

- Molecule 3 is water.

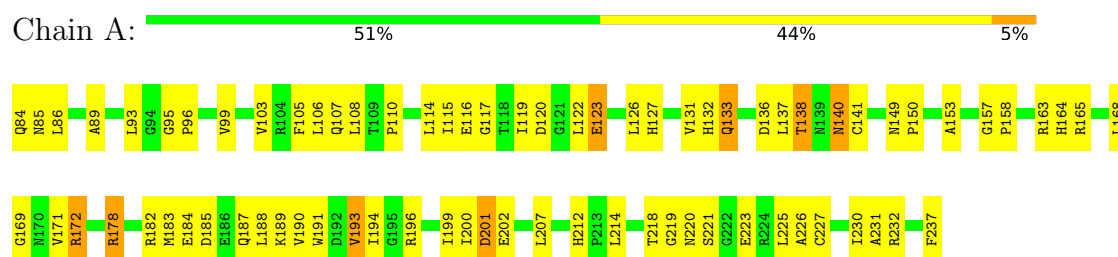
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	6	Total	O	0	0
			6	6		
3	B	7	Total	O	0	0
			7	7		
3	C	8	Total	O	0	0
			8	8		
3	D	5	Total	O	0	0
			5	5		

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. 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.

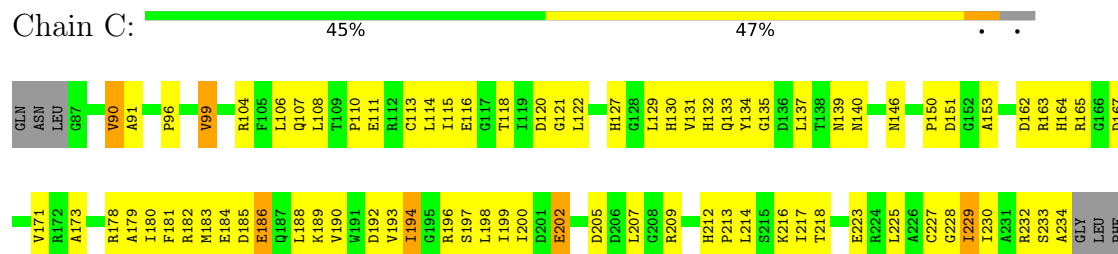
• Molecule 1: HUMAN COPPER CHAPERONE FOR SUPEROXIDE DISMUTASE DOMAIN II



• Molecule 1: HUMAN COPPER CHAPERONE FOR SUPEROXIDE DISMUTASE DOMAIN II



• Molecule 1: HUMAN COPPER CHAPERONE FOR SUPEROXIDE DISMUTASE DOMAIN II



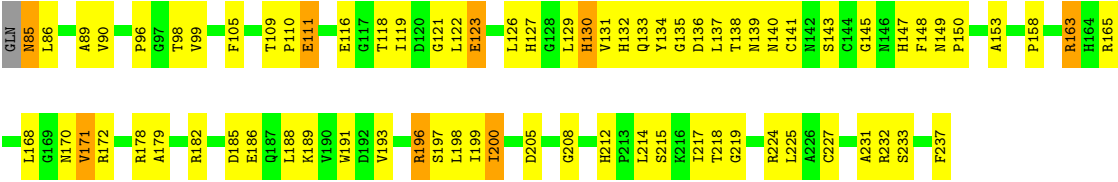
• Molecule 1: HUMAN COPPER CHAPERONE FOR SUPEROXIDE DISMUTASE DOMAIN II

Chain D:

51%

44%

5% •



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	117.19Å 66.72Å 88.03Å 90.00° 96.63° 90.00°	Depositor
Resolution (Å)	19.21 – 2.75	Depositor
% Data completeness (in resolution range)	90.1 (19.21-2.75)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.228 , 0.283	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4495	wwPDB-VP
Average B, all atoms (Å ²)	32.0	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: ZN

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	0.47	0/1162	1.06	8/1568 (0.5%)
1	B	0.45	0/1112	1.00	5/1502 (0.3%)
1	C	0.53	1/1112 (0.1%)	0.99	0/1502
1	D	0.45	0/1153	0.97	2/1556 (0.1%)
All	All	0.48	1/4539 (0.0%)	1.01	15/6128 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	216	LYS	C-N	-7.42	1.25	1.33

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	95	GLY	CA-C-N	7.62	127.79	119.87
1	A	95	GLY	C-N-CA	7.62	127.79	119.87
1	A	219	GLY	N-CA-C	-6.85	106.25	115.36
1	B	149	ASN	CA-C-N	6.43	125.66	118.97
1	B	149	ASN	C-N-CA	6.43	125.66	118.97
1	D	219	GLY	N-CA-C	-6.36	106.30	115.27
1	B	219	GLY	N-CA-C	-6.21	106.58	115.64
1	D	123	GLU	N-CA-C	-6.01	102.08	109.65
1	A	172	ARG	N-CA-C	5.71	117.60	108.41
1	B	193	VAL	N-CA-C	5.59	117.58	112.43
1	A	201	ASP	N-CA-C	5.55	118.74	110.52
1	A	133	GLN	N-CA-C	5.54	118.25	111.82
1	B	170	ASN	N-CA-C	5.34	118.36	110.46
1	A	123	GLU	N-CA-C	-5.09	103.61	109.93
1	A	193	VAL	N-CA-C	5.02	119.53	113.00

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	1143	0	1093	76	0
1	B	1094	0	1045	73	0
1	C	1094	0	1045	86	0
1	D	1134	0	1085	61	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	6	0	0	0	0
3	B	7	0	0	1	0
3	C	8	0	0	1	0
3	D	5	0	0	1	0
All	All	4495	0	4268	278	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 32.

All (278) 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:150:PRO:HD2	1:B:165:ARG:HG3	1.44	0.94
1:A:212:HIS:HD2	1:A:214:LEU:H	1.17	0.88
1:C:133:GLN:HE21	1:C:146:ASN:H	0.90	0.88
1:D:149:ASN:HD21	1:D:153:ALA:H	1.21	0.87
1:C:90:VAL:HG23	1:C:233:SER:HB2	1.61	0.82
1:D:90:VAL:HG23	1:D:233:SER:HB2	1.59	0.82
1:C:133:GLN:NE2	1:C:146:ASN:H	1.74	0.81
1:D:130:HIS:HE1	1:D:205:ASP:OD2	1.61	0.81
1:B:119:ILE:HG21	1:B:200:ILE:HD13	1.60	0.81
1:B:168:LEU:HD21	1:B:188:LEU:HD22	1.63	0.81
1:B:109:THR:HB	1:B:110:PRO:HD2	1.61	0.80

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:133:GLN:HE21	1:C:146:ASN:N	1.75	0.80
1:B:119:ILE:HG21	1:B:200:ILE:CD1	2.12	0.79
1:D:85:ASN:N	1:D:85:ASN:HD22	1.79	0.79
1:A:199:ILE:HG12	1:A:227:CYS:HB3	1.65	0.78
1:D:135:GLY:HA2	1:D:197:SER:HB3	1.66	0.77
1:A:212:HIS:CD2	1:A:214:LEU:H	2.00	0.77
1:A:150:PRO:HD2	1:A:165:ARG:HG3	1.66	0.77
1:D:109:THR:HB	1:D:110:PRO:HD2	1.69	0.74
1:A:99:VAL:HG21	1:A:225:LEU:HD13	1.70	0.73
1:A:149:ASN:HD21	1:A:153:ALA:H	1.38	0.72
1:D:99:VAL:HG11	1:D:200:ILE:HD12	1.71	0.71
1:C:188:LEU:HD11	1:C:193:VAL:HG11	1.73	0.70
1:D:89:ALA:HB3	1:D:105:PHE:HB2	1.75	0.68
1:D:217:ILE:HG22	1:D:218:THR:HG23	1.75	0.68
1:A:199:ILE:HG12	1:A:227:CYS:CB	2.24	0.68
1:D:123:GLU:O	1:D:127:HIS:HE1	1.77	0.68
1:C:212:HIS:CD2	1:C:214:LEU:H	2.12	0.67
1:D:111:GLU:HA	1:D:189:LYS:HE3	1.76	0.67
1:C:202:GLU:HA	1:C:225:LEU:HD11	1.78	0.66
1:A:136:ASP:OD1	1:A:138:THR:HB	1.96	0.66
1:C:212:HIS:HD2	1:C:214:LEU:H	1.44	0.65
1:B:130:HIS:HE1	1:B:205:ASP:OD2	1.78	0.65
1:D:85:ASN:N	1:D:85:ASN:ND2	2.45	0.65
1:C:153:ALA:HB1	1:C:162:ASP:OD1	1.97	0.65
1:B:200:ILE:HG13	1:B:226:ALA:HB3	1.79	0.65
1:C:114:LEU:HD21	1:C:116:GLU:OE2	1.98	0.64
1:C:171:VAL:HG12	1:C:179:ALA:HB1	1.80	0.64
1:C:107:GLN:NE2	1:C:189:LYS:HD2	2.13	0.64
1:C:199:ILE:HG12	1:C:227:CYS:HB3	1.80	0.64
1:D:150:PRO:HG2	1:D:165:ARG:NH1	2.12	0.64
1:D:208:GLY:HA2	1:D:215:SER:O	1.98	0.64
1:B:99:VAL:HG11	1:B:225:LEU:HD13	1.79	0.63
1:A:193:VAL:HA	1:A:196:ARG:HD2	1.80	0.63
1:A:194:ILE:HD13	1:A:232:ARG:HG2	1.80	0.63
1:D:96:PRO:HD2	1:D:225:LEU:O	1.99	0.62
1:B:183:MET:HE2	1:B:185:ASP:HB2	1.80	0.62
1:C:122:LEU:HD21	1:C:200:ILE:HD13	1.81	0.62
1:A:232:ARG:HD2	1:B:196:ARG:HG2	1.81	0.62
1:B:178:ARG:NE	1:B:180:ILE:HD11	2.15	0.62
1:B:190:VAL:O	1:B:194:ILE:HG23	2.00	0.62
1:D:149:ASN:HD21	1:D:153:ALA:N	1.96	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:178:ARG:CZ	1:B:180:ILE:HD11	2.30	0.62
1:C:115:ILE:HB	1:C:183:MET:HG3	1.82	0.61
1:A:108:LEU:HD12	1:A:184:GLU:OE2	2.00	0.61
1:B:138:THR:HG22	1:C:139:ASN:OD1	2.01	0.61
1:A:123:GLU:O	1:A:127:HIS:HE1	1.83	0.61
1:C:189:LYS:O	1:C:192:ASP:HB2	2.01	0.61
1:D:178:ARG:HD2	3:D:23:HOH:O	2.00	0.60
1:B:157:GLY:H	1:B:160:ASP:HB2	1.65	0.60
1:A:237:PHE:O	1:C:146:ASN:HB3	2.01	0.60
1:D:119:ILE:HG21	1:D:200:ILE:HD13	1.82	0.60
1:A:194:ILE:HG23	1:B:195:GLY:HA3	1.84	0.60
1:C:133:GLN:HG3	1:C:134:TYR:CD1	2.36	0.59
1:B:188:LEU:HD11	1:B:193:VAL:HG11	1.83	0.59
1:C:188:LEU:HD11	1:C:193:VAL:CG1	2.31	0.59
1:A:116:GLU:HG2	1:A:182:ARG:HG3	1.85	0.59
1:A:188:LEU:HD11	1:A:193:VAL:CG1	2.33	0.59
1:B:108:LEU:HD11	1:B:114:LEU:HB2	1.85	0.59
1:A:232:ARG:O	1:B:134:TYR:HB3	2.03	0.58
1:C:122:LEU:HB3	1:C:127:HIS:CE1	2.39	0.58
1:B:139:ASN:O	1:B:140:ASN:HB2	2.02	0.58
1:B:163:ARG:NH2	1:B:187:GLN:HB3	2.19	0.58
1:B:199:ILE:HG12	1:B:227:CYS:CB	2.34	0.58
1:C:196:ARG:HG2	1:D:232:ARG:HD2	1.84	0.58
1:A:149:ASN:HD21	1:A:153:ALA:N	2.00	0.57
1:C:193:VAL:HB	1:C:230:ILE:HD12	1.86	0.57
1:A:157:GLY:HA2	1:A:207:LEU:HD22	1.85	0.57
1:D:116:GLU:HG2	1:D:182:ARG:HG3	1.86	0.57
1:A:200:ILE:HG23	1:A:225:LEU:HD12	1.85	0.57
1:B:149:ASN:HD21	1:B:153:ALA:H	1.53	0.57
1:B:189:LYS:HG3	1:B:191:TRP:CZ2	2.39	0.57
1:A:133:GLN:O	1:A:196:ARG:NH1	2.37	0.57
1:C:115:ILE:HD11	1:C:188:LEU:HD23	1.87	0.57
1:C:228:GLY:O	1:C:229:ILE:HB	2.05	0.57
1:B:139:ASN:ND2	3:B:13:HOH:O	2.38	0.56
1:C:99:VAL:HA	1:C:120:ASP:O	2.05	0.56
1:D:145:GLY:O	1:D:224:ARG:NH2	2.35	0.56
1:A:103:VAL:HG22	1:A:117:GLY:HA3	1.87	0.56
1:C:217:ILE:HG22	1:C:218:THR:HG23	1.88	0.56
1:A:132:HIS:O	1:A:196:ARG:HB3	2.06	0.56
1:B:139:ASN:CG	1:C:139:ASN:HB2	2.30	0.56
1:C:114:LEU:HD21	1:C:116:GLU:CD	2.30	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:135:GLY:HA2	1:C:197:SER:HB3	1.88	0.56
1:B:124:PRO:HA	1:B:173:ALA:O	2.06	0.56
1:D:99:VAL:CG1	1:D:200:ILE:HD12	2.35	0.55
1:B:194:ILE:HD12	1:B:232:ARG:CZ	2.36	0.55
1:C:110:PRO:HG2	1:C:111:GLU:OE2	2.06	0.55
1:A:127:HIS:HD2	1:A:201:ASP:O	1.90	0.55
1:B:107:GLN:OE1	1:B:113:CYS:HB2	2.07	0.55
1:B:136:ASP:OD1	1:B:136:ASP:C	2.50	0.55
1:C:130:HIS:HE1	1:C:205:ASP:OD2	1.90	0.54
1:D:136:ASP:OD1	1:D:138:THR:HB	2.07	0.54
1:C:132:HIS:CE1	1:C:199:ILE:HD12	2.42	0.54
1:B:110:PRO:C	1:B:111:GLU:HG3	2.33	0.54
1:B:120:ASP:OD1	1:B:178:ARG:HG3	2.08	0.54
1:D:132:HIS:O	1:D:196:ARG:HB3	2.08	0.54
1:D:158:PRO:HG3	1:D:168:LEU:C	2.33	0.53
1:A:158:PRO:HG3	1:A:168:LEU:C	2.34	0.53
1:C:228:GLY:O	1:C:229:ILE:CB	2.56	0.53
1:B:164:HIS:HB2	1:B:167:ASP:CG	2.32	0.53
1:B:122:LEU:HD21	1:B:200:ILE:HD12	1.91	0.53
1:A:212:HIS:HD2	1:A:214:LEU:N	1.98	0.53
1:B:90:VAL:HG21	1:B:104:ARG:NH2	2.24	0.53
1:B:99:VAL:HA	1:B:120:ASP:O	2.08	0.53
1:D:212:HIS:HD2	1:D:214:LEU:HB2	1.73	0.53
1:A:132:HIS:CE1	1:A:199:ILE:HD12	2.44	0.53
1:B:151:ASP:OD2	1:B:165:ARG:HD3	2.09	0.53
1:B:199:ILE:HG12	1:B:227:CYS:HB3	1.91	0.53
1:D:119:ILE:HG21	1:D:200:ILE:CD1	2.39	0.52
1:D:132:HIS:CE1	1:D:199:ILE:HD12	2.44	0.52
1:C:190:VAL:O	1:C:194:ILE:HG23	2.10	0.52
1:B:126:LEU:HD21	1:B:172:ARG:NH1	2.24	0.52
1:C:163:ARG:HH22	1:C:185:ASP:CG	2.17	0.52
1:C:193:VAL:HG23	1:C:230:ILE:HG21	1.92	0.52
1:B:99:VAL:HG11	1:B:225:LEU:CD1	2.39	0.52
1:A:110:PRO:O	1:A:189:LYS:HE3	2.10	0.52
1:D:137:LEU:HA	1:D:140:ASN:O	2.09	0.51
1:A:107:GLN:HB2	1:A:190:VAL:CG2	2.41	0.51
1:A:165:ARG:HD2	1:A:187:GLN:HE21	1.76	0.51
1:D:189:LYS:HD2	1:D:191:TRP:CH2	2.45	0.51
1:A:96:PRO:HD2	1:A:225:LEU:O	2.11	0.51
1:A:188:LEU:HD11	1:A:193:VAL:HG11	1.91	0.51
1:D:199:ILE:HG12	1:D:227:CYS:HB3	1.93	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:202:GLU:HB2	1:C:225:LEU:HD21	1.93	0.51
1:D:141:CYS:O	1:D:224:ARG:NH1	2.42	0.50
1:B:101:GLY:HA3	1:B:118:THR:O	2.11	0.50
1:A:193:VAL:HB	1:A:230:ILE:CD1	2.41	0.50
1:C:150:PRO:CG	1:C:165:ARG:HG3	2.40	0.50
1:B:138:THR:CG2	3:C:16:HOH:O	2.59	0.50
1:B:123:GLU:O	1:B:124:PRO:C	2.55	0.50
1:A:165:ARG:CD	1:A:187:GLN:HE21	2.24	0.50
1:B:109:THR:OG1	1:B:112:ARG:HG2	2.12	0.50
1:C:194:ILE:HD13	1:C:232:ARG:CZ	2.42	0.50
1:D:99:VAL:HG13	1:D:122:LEU:HG	1.94	0.50
1:A:89:ALA:HA	1:A:231:ALA:O	2.12	0.49
1:C:173:ALA:HA	1:C:179:ALA:HB2	1.93	0.49
1:C:99:VAL:H	1:C:121:GLY:HA3	1.77	0.49
1:C:116:GLU:OE2	1:C:182:ARG:CZ	2.61	0.49
1:D:193:VAL:HA	1:D:196:ARG:HD2	1.94	0.49
1:B:123:GLU:O	1:B:127:HIS:HE1	1.95	0.49
1:A:200:ILE:CG2	1:A:226:ALA:HB3	2.42	0.49
1:D:126:LEU:HD23	1:D:172:ARG:HB2	1.94	0.49
1:D:163:ARG:HH22	1:D:185:ASP:CG	2.20	0.49
1:B:93:LEU:HD23	1:B:228:GLY:N	2.28	0.48
1:B:158:PRO:HG3	1:B:168:LEU:C	2.37	0.48
1:B:188:LEU:CD1	1:B:193:VAL:HG11	2.43	0.48
1:B:208:GLY:HA2	1:B:215:SER:O	2.13	0.48
1:C:139:ASN:O	1:C:140:ASN:HB2	2.13	0.48
1:A:193:VAL:HG23	1:A:230:ILE:HG21	1.95	0.48
1:A:196:ARG:HG2	1:B:232:ARG:HD2	1.94	0.48
1:A:202:GLU:N	1:A:223:GLU:O	2.45	0.48
1:A:106:LEU:O	1:A:114:LEU:N	2.45	0.48
1:A:150:PRO:HD2	1:A:165:ARG:CG	2.40	0.48
1:D:147:HIS:HE1	1:D:218:THR:O	1.96	0.48
1:C:131:VAL:HA	1:C:197:SER:O	2.14	0.48
1:C:171:VAL:CG1	1:C:179:ALA:HB1	2.42	0.48
1:A:108:LEU:HB3	1:C:213:PRO:CB	2.44	0.47
1:A:165:ARG:HA	1:A:187:GLN:NE2	2.29	0.47
1:B:90:VAL:HG21	1:B:104:ARG:HH21	1.80	0.47
1:C:118:THR:OG1	1:C:180:ILE:HG12	2.14	0.47
1:C:182:ARG:HG2	1:C:182:ARG:HH11	1.80	0.47
1:C:193:VAL:HA	1:C:196:ARG:HD2	1.95	0.47
1:C:212:HIS:HD2	1:C:214:LEU:HB2	1.80	0.47
1:A:99:VAL:CG2	1:A:225:LEU:HD13	2.42	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:131:VAL:HG21	1:D:188:LEU:HD21	1.96	0.47
1:A:140:ASN:HD22	1:A:140:ASN:HA	1.56	0.47
1:A:158:PRO:CG	1:A:169:GLY:HA3	2.45	0.47
1:C:137:LEU:HD11	1:C:229:ILE:HG13	1.96	0.46
1:A:158:PRO:HG3	1:A:169:GLY:N	2.30	0.46
1:C:133:GLN:O	1:C:196:ARG:HG2	2.16	0.46
1:A:120:ASP:OD1	1:A:178:ARG:HB2	2.15	0.46
1:B:116:GLU:HG2	1:B:182:ARG:HE	1.80	0.46
1:D:199:ILE:HG12	1:D:227:CYS:CB	2.46	0.46
1:B:218:THR:CG2	1:D:237:PHE:HA	2.46	0.46
1:C:115:ILE:HG22	1:C:181:PHE:HE1	1.81	0.46
1:A:165:ARG:HD2	1:A:187:GLN:HG2	1.98	0.46
1:A:201:ASP:HB3	1:A:221:SER:O	2.15	0.45
1:B:218:THR:HG21	1:D:237:PHE:HA	1.99	0.45
1:C:90:VAL:HG12	1:C:91:ALA:H	1.81	0.45
1:D:86:LEU:N	1:D:86:LEU:HD22	2.32	0.45
1:D:89:ALA:HA	1:D:231:ALA:O	2.16	0.45
1:B:94:GLY:O	1:B:226:ALA:HA	2.16	0.45
1:B:105:PHE:CD2	1:B:115:ILE:HG12	2.51	0.45
1:C:90:VAL:HG21	1:D:136:ASP:OD2	2.15	0.45
1:D:133:GLN:HE21	1:D:145:GLY:HA3	1.82	0.45
1:D:133:GLN:O	1:D:196:ARG:NH1	2.48	0.45
1:C:202:GLU:HA	1:C:225:LEU:CD1	2.43	0.45
1:A:131:VAL:HG11	1:A:193:VAL:HG12	1.98	0.45
1:A:218:THR:C	1:A:220:ASN:H	2.24	0.45
1:C:90:VAL:HG12	1:C:91:ALA:N	2.31	0.45
1:D:110:PRO:O	1:D:189:LYS:HE3	2.17	0.45
1:D:130:HIS:CE1	1:D:205:ASP:OD2	2.54	0.45
1:C:150:PRO:HG2	1:C:165:ARG:HG3	1.98	0.45
1:C:107:GLN:NE2	1:C:113:CYS:HB2	2.32	0.45
1:D:129:LEU:HD22	1:D:171:VAL:HG21	1.99	0.45
1:B:131:VAL:HA	1:B:197:SER:O	2.18	0.44
1:B:217:ILE:HD11	1:D:86:LEU:HD21	1.99	0.44
1:C:129:LEU:HD11	1:C:198:LEU:HD11	1.99	0.44
1:D:123:GLU:O	1:D:127:HIS:CE1	2.66	0.44
1:A:149:ASN:CG	1:A:149:ASN:O	2.57	0.44
1:C:113:CYS:O	1:C:184:GLU:HA	2.17	0.44
1:A:93:LEU:HD12	1:A:119:ILE:HD13	1.99	0.44
1:C:129:LEU:HD22	1:C:171:VAL:HG21	1.99	0.44
1:A:105:PHE:CD2	1:A:115:ILE:HG12	2.53	0.43
1:B:198:LEU:O	1:B:227:CYS:HA	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:118:THR:HG21	1:C:178:ARG:HE	1.84	0.43
1:A:122:LEU:HD21	1:A:200:ILE:HG12	1.99	0.43
1:A:172:ARG:NH1	1:A:172:ARG:HG2	2.33	0.43
1:C:96:PRO:HD2	1:C:225:LEU:O	2.18	0.43
1:C:186:GLU:H	1:C:186:GLU:HG2	1.67	0.43
1:D:118:THR:HA	1:D:179:ALA:O	2.18	0.43
1:D:130:HIS:O	1:D:198:LEU:HD12	2.18	0.43
1:B:148:PHE:CE1	1:B:196:ARG:CZ	3.02	0.43
1:C:122:LEU:HD11	1:C:200:ILE:CD1	2.49	0.43
1:A:172:ARG:HG2	1:A:172:ARG:HH11	1.84	0.43
1:B:212:HIS:O	1:B:213:PRO:C	2.62	0.43
1:C:234:ALA:HB2	1:D:134:TYR:CZ	2.54	0.43
1:B:148:PHE:CE1	1:B:150:PRO:HG3	2.53	0.43
1:B:164:HIS:HB2	1:B:167:ASP:OD2	2.19	0.43
1:A:140:ASN:HB3	1:A:141:CYS:H	1.60	0.43
1:A:189:LYS:HD2	1:A:191:TRP:CZ2	2.54	0.43
1:B:139:ASN:HB2	1:C:139:ASN:HB2	2.01	0.43
1:D:186:GLU:O	1:D:189:LYS:NZ	2.52	0.42
1:B:113:CYS:O	1:B:184:GLU:HA	2.18	0.42
1:C:209:ARG:HH11	1:C:209:ARG:HG2	1.84	0.42
1:A:200:ILE:HG23	1:A:225:LEU:CD1	2.48	0.42
1:D:98:THR:HG22	1:D:98:THR:O	2.19	0.42
1:D:99:VAL:HG22	1:D:121:GLY:O	2.19	0.42
1:A:99:VAL:HA	1:A:120:ASP:O	2.19	0.42
1:B:163:ARG:HD2	1:B:164:HIS:O	2.20	0.42
1:C:137:LEU:HD22	1:C:227:CYS:SG	2.59	0.42
1:C:205:ASP:OD1	1:C:207:LEU:N	2.48	0.42
1:A:108:LEU:HB3	1:C:213:PRO:HB2	2.01	0.42
1:D:186:GLU:H	1:D:186:GLU:HG2	1.68	0.42
1:C:205:ASP:OD1	1:C:207:LEU:HD23	2.19	0.42
1:C:90:VAL:CG2	1:C:233:SER:HB2	2.42	0.42
1:A:196:ARG:HG2	1:A:196:ARG:HH11	1.84	0.42
1:B:199:ILE:HG12	1:B:227:CYS:HB2	2.01	0.42
1:A:194:ILE:CG2	1:B:195:GLY:HA3	2.48	0.41
1:A:137:LEU:HD22	1:A:227:CYS:SG	2.60	0.41
1:B:98:THR:O	1:B:99:VAL:HB	2.19	0.41
1:A:86:LEU:HD11	1:C:217:ILE:HD12	2.02	0.41
1:B:197:SER:HA	1:B:230:ILE:HG13	2.02	0.41
1:D:126:LEU:HD22	1:D:170:ASN:HB3	2.02	0.41
1:C:107:GLN:CD	1:C:113:CYS:HB2	2.45	0.41
1:C:207:LEU:O	1:C:209:ARG:HG3	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:LEU:HD23	1:A:172:ARG:HB2	2.03	0.41
1:A:84:GLN:NE2	1:A:85:ASN:H	2.18	0.41
1:A:196:ARG:NH1	1:A:196:ARG:HG2	2.36	0.41
1:C:133:GLN:HG3	1:C:134:TYR:CE1	2.55	0.41
1:C:106:LEU:HG	1:C:108:LEU:HD23	2.02	0.41
1:D:148:PHE:CZ	1:D:150:PRO:HG3	2.55	0.41
1:A:107:GLN:HB2	1:A:190:VAL:HG21	2.02	0.41
1:A:163:ARG:CD	1:A:164:HIS:O	2.69	0.41
1:A:163:ARG:HH22	1:A:185:ASP:CG	2.29	0.41
1:A:168:LEU:HD21	1:A:188:LEU:HD22	2.02	0.41
1:B:90:VAL:HG23	1:B:233:SER:HB2	2.03	0.41
1:C:164:HIS:HB2	1:C:167:ASP:CG	2.46	0.41
1:C:131:VAL:HG22	1:C:230:ILE:HD11	2.03	0.40
1:C:135:GLY:O	1:C:229:ILE:HG12	2.21	0.40
1:A:126:LEU:CD2	1:A:172:ARG:HB2	2.51	0.40
1:B:99:VAL:H	1:B:121:GLY:HA3	1.85	0.40
1:B:174:ASP:C	1:B:174:ASP:OD2	2.64	0.40
1:C:151:ASP:OD1	1:C:165:ARG:NH1	2.55	0.40
1:C:90:VAL:HG13	1:C:104:ARG:HG2	2.03	0.40
1:D:139:ASN:O	1:D:140:ASN:HB2	2.21	0.40
1:C:114:LEU:HG	1:C:116:GLU:HG3	2.03	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	152/154 (99%)	141 (93%)	11 (7%)	0	100	100
1	B	146/154 (95%)	126 (86%)	16 (11%)	4 (3%)	4	7
1	C	146/154 (95%)	130 (89%)	14 (10%)	2 (1%)	9	17
1	D	151/154 (98%)	145 (96%)	5 (3%)	1 (1%)	18	34

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	595/616 (97%)	542 (91%)	46 (8%)	7 (1%)	10	20

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	111	GLU
1	C	229	ILE
1	D	196	ARG
1	B	99	VAL
1	B	110	PRO
1	B	140	ASN
1	C	99	VAL

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	118/118 (100%)	113 (96%)	5 (4%)	26	49
1	B	113/118 (96%)	108 (96%)	5 (4%)	25	47
1	C	113/118 (96%)	108 (96%)	5 (4%)	25	47
1	D	117/118 (99%)	110 (94%)	7 (6%)	17	32
All	All	461/472 (98%)	439 (95%)	22 (5%)	23	44

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	138	THR
1	A	140	ASN
1	A	171	VAL
1	A	178	ARG
1	A	183	MET
1	B	111	GLU
1	B	143	SER
1	B	165	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	171	VAL
1	B	200	ILE
1	C	90	VAL
1	C	186	GLU
1	C	194	ILE
1	C	202	GLU
1	C	223	GLU
1	D	85	ASN
1	D	111	GLU
1	D	130	HIS
1	D	143	SER
1	D	163	ARG
1	D	171	VAL
1	D	200	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	84	GLN
1	A	127	HIS
1	A	132	HIS
1	A	139	ASN
1	A	140	ASN
1	A	142	ASN
1	A	149	ASN
1	A	187	GLN
1	A	212	HIS
1	B	127	HIS
1	B	130	HIS
1	B	139	ASN
1	B	140	ASN
1	B	142	ASN
1	B	149	ASN
1	B	164	HIS
1	C	130	HIS
1	C	133	GLN
1	C	140	ASN
1	C	142	ASN
1	C	149	ASN
1	C	212	HIS
1	D	127	HIS
1	D	130	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	132	HIS
1	D	133	GLN
1	D	139	ASN
1	D	140	ASN
1	D	142	ASN
1	D	146	ASN
1	D	149	ASN
1	D	187	GLN
1	D	212	HIS

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 4 ligands modelled in this entry, 4 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 ⓘ

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.