



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:35 AM UTC

PDB ID : 4OUL / pdb_00004oul
Title : Crystal structure of human Caprin-2 C1q domain
Authors : Song, X.; Li, L.
Deposited on : 2014-02-17
Resolution : 1.95 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
Mogul : 2022.3.0, CSD as543be (2022)
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

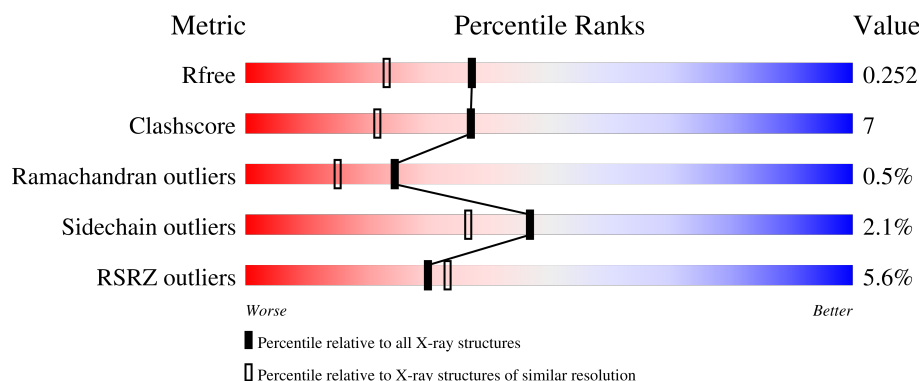
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	140	8% 74% 19% • 6%
1	B	140	4% 79% 14% • 6%
1	C	140	3% 81% 12% 6%
1	D	140	6% 79% 14% • 6%
1	E	140	4% 76% 16% • 7%

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Mol	Chain	Length	Quality of chain
1	F	140	76% 18% 6% 7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6291 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 Caprin-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	132	Total	C	N	O	S	0	1	0
			1042	675	178	186	3			
1	B	132	Total	C	N	O	S	0	0	0
			1010	653	168	186	3			
1	C	131	Total	C	N	O	S	0	0	0
			1003	653	166	181	3			
1	D	132	Total	C	N	O	S	0	0	0
			1024	664	169	188	3			
1	E	130	Total	C	N	O	S	0	0	0
			1017	660	170	184	3			
1	F	132	Total	C	N	O	S	0	0	0
			1043	674	175	190	4			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1128	LEU	-	expression tag	UNP Q6IMN6
A	1129	GLU	-	expression tag	UNP Q6IMN6
A	1130	HIS	-	expression tag	UNP Q6IMN6
A	1131	HIS	-	expression tag	UNP Q6IMN6
A	1132	HIS	-	expression tag	UNP Q6IMN6
A	1133	HIS	-	expression tag	UNP Q6IMN6
A	1134	HIS	-	expression tag	UNP Q6IMN6
A	1135	HIS	-	expression tag	UNP Q6IMN6
B	1128	LEU	-	expression tag	UNP Q6IMN6
B	1129	GLU	-	expression tag	UNP Q6IMN6
B	1130	HIS	-	expression tag	UNP Q6IMN6
B	1131	HIS	-	expression tag	UNP Q6IMN6
B	1132	HIS	-	expression tag	UNP Q6IMN6
B	1133	HIS	-	expression tag	UNP Q6IMN6
B	1134	HIS	-	expression tag	UNP Q6IMN6
B	1135	HIS	-	expression tag	UNP Q6IMN6
C	1128	LEU	-	expression tag	UNP Q6IMN6

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1129	GLU	-	expression tag	UNP Q6IMN6
C	1130	HIS	-	expression tag	UNP Q6IMN6
C	1131	HIS	-	expression tag	UNP Q6IMN6
C	1132	HIS	-	expression tag	UNP Q6IMN6
C	1133	HIS	-	expression tag	UNP Q6IMN6
C	1134	HIS	-	expression tag	UNP Q6IMN6
C	1135	HIS	-	expression tag	UNP Q6IMN6
D	1128	LEU	-	expression tag	UNP Q6IMN6
D	1129	GLU	-	expression tag	UNP Q6IMN6
D	1130	HIS	-	expression tag	UNP Q6IMN6
D	1131	HIS	-	expression tag	UNP Q6IMN6
D	1132	HIS	-	expression tag	UNP Q6IMN6
D	1133	HIS	-	expression tag	UNP Q6IMN6
D	1134	HIS	-	expression tag	UNP Q6IMN6
D	1135	HIS	-	expression tag	UNP Q6IMN6
E	1128	LEU	-	expression tag	UNP Q6IMN6
E	1129	GLU	-	expression tag	UNP Q6IMN6
E	1130	HIS	-	expression tag	UNP Q6IMN6
E	1131	HIS	-	expression tag	UNP Q6IMN6
E	1132	HIS	-	expression tag	UNP Q6IMN6
E	1133	HIS	-	expression tag	UNP Q6IMN6
E	1134	HIS	-	expression tag	UNP Q6IMN6
E	1135	HIS	-	expression tag	UNP Q6IMN6
F	1128	LEU	-	expression tag	UNP Q6IMN6
F	1129	GLU	-	expression tag	UNP Q6IMN6
F	1130	HIS	-	expression tag	UNP Q6IMN6
F	1131	HIS	-	expression tag	UNP Q6IMN6
F	1132	HIS	-	expression tag	UNP Q6IMN6
F	1133	HIS	-	expression tag	UNP Q6IMN6
F	1134	HIS	-	expression tag	UNP Q6IMN6
F	1135	HIS	-	expression tag	UNP Q6IMN6

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Ca 1 1	0	0
2	B	1	Total Ca 1 1	0	0
2	D	1	Total Ca 1 1	0	0
2	E	1	Total Ca 1 1	0	0

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			6	3	3		

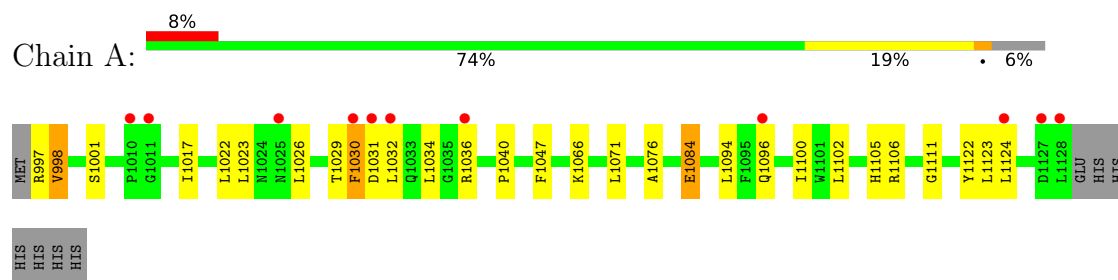
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	21	Total	O	0	0
			21	21		
4	B	28	Total	O	0	0
			28	28		
4	C	23	Total	O	0	0
			23	23		
4	D	24	Total	O	0	0
			24	24		
4	E	28	Total	O	0	0
			28	28		
4	F	18	Total	O	0	0
			18	18		

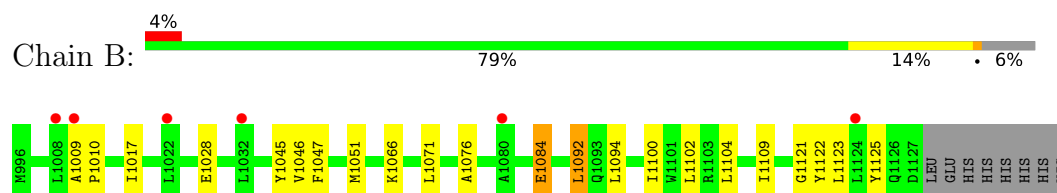
3 Residue-property plots [i](#)

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.

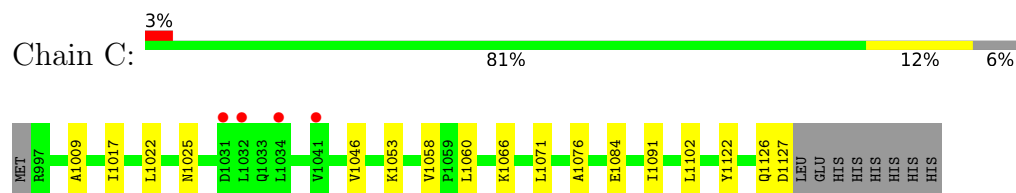
- Molecule 1: Caprin-2



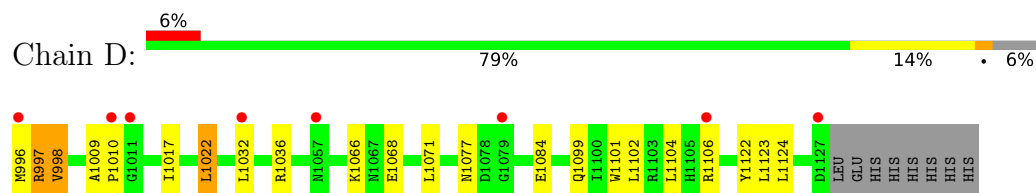
- Molecule 1: Caprin-2



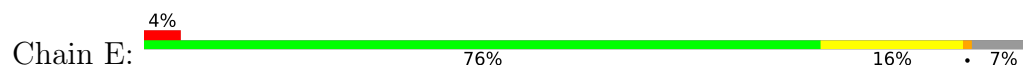
- Molecule 1: Caprin-2



- Molecule 1: Caprin-2

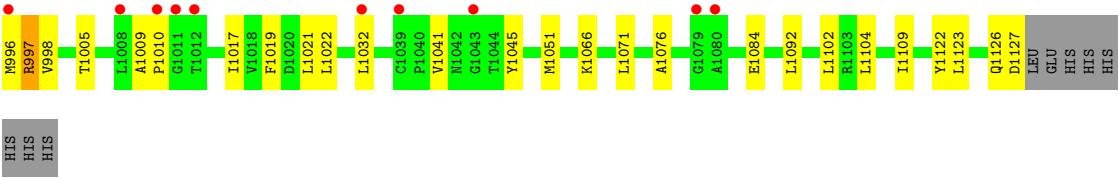
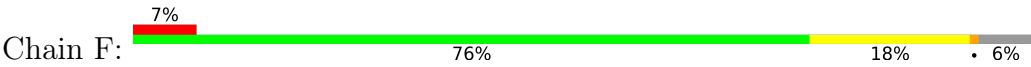


- Molecule 1: Caprin-2





● Molecule 1: Caprin-2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	48.35Å 48.34Å 80.88Å 89.51° 87.24° 71.74°	Depositor
Resolution (Å)	45.90 – 1.95 45.90 – 1.95	Depositor EDS
% Data completeness (in resolution range)	89.8 (45.90-1.95) 84.1 (45.90-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 1.95Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.203 , 0.251 0.208 , 0.252	Depositor DCC
R_{free} test set	1979 reflections (3.90%)	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.330	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.015 for -h,-k,l 0.007 for k,h,-l 0.015 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6291	wwPDB-VP
Average B, all atoms (Å ²)	36.0	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 40.37 % 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 2.7616e-04. 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

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

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	0.56	0/1069	0.86	1/1457 (0.1%)
1	B	0.50	0/1036	0.82	2/1416 (0.1%)
1	C	0.49	0/1030	0.88	2/1407 (0.1%)
1	D	0.53	0/1051	0.87	3/1433 (0.2%)
1	E	0.54	0/1043	0.82	0/1420
1	F	0.50	0/1069	0.93	3/1457 (0.2%)
All	All	0.52	0/6298	0.86	11/8590 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	1009	ALA	CA-C-N	8.55	128.28	119.56
1	F	1009	ALA	C-N-CA	8.55	128.28	119.56
1	C	1009	ALA	CA-C-N	7.09	126.92	119.05
1	C	1009	ALA	C-N-CA	7.09	126.92	119.05
1	D	1009	ALA	CA-C-N	7.07	127.06	119.28
1	D	1009	ALA	C-N-CA	7.07	127.06	119.28
1	A	1029	THR	N-CA-C	-6.91	99.74	110.17
1	F	998	VAL	N-CA-C	6.57	117.49	107.77
1	D	998	VAL	N-CA-C	6.08	116.77	107.77
1	B	1009	ALA	CA-C-N	5.27	125.21	119.78
1	B	1009	ALA	C-N-CA	5.27	125.21	119.78

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	1042	0	1009	22	0
1	B	1010	0	944	14	0
1	C	1003	0	947	10	0
1	D	1024	0	973	15	0
1	E	1017	0	978	15	0
1	F	1043	0	1009	20	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
3	D	6	0	8	0	0
4	A	21	0	0	1	0
4	B	28	0	0	0	0
4	C	23	0	0	1	0
4	D	24	0	0	1	0
4	E	28	0	0	2	0
4	F	18	0	0	1	0
All	All	6291	0	5868	85	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 7.

All (85) 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:C:1076:ALA:HB1	1:C:1084:GLU:HG3	1.56	0.86
1:E:1066:LYS:HB2	1:E:1071:LEU:HD11	1.62	0.81
1:E:1076:ALA:HB1	1:E:1084:GLU:HG3	1.62	0.81
1:D:1066:LYS:HB2	1:D:1071:LEU:HD11	1.63	0.81
1:A:1040:PRO:O	1:A:1096:GLN:NE2	2.19	0.76
1:F:997:ARG:NH1	1:F:1126:GLN:OE1	2.24	0.70
1:D:1077:ASN:OD1	4:D:1320:HOH:O	2.08	0.70
1:E:1087:SER:OG	4:E:1320:HOH:O	2.10	0.69
1:F:1066:LYS:HB2	1:F:1071:LEU:HD11	1.73	0.69
1:A:1066:LYS:HB2	1:A:1071:LEU:HD11	1.75	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1066:LYS:HB2	1:C:1071:LEU:HD11	1.75	0.68
1:F:996:MET:O	1:F:997:ARG:HB2	1.92	0.68
1:C:1025:ASN:OD1	4:C:1214:HOH:O	2.12	0.67
1:D:1122:TYR:OH	1:F:1122:TYR:OH	2.12	0.66
1:D:996:MET:O	1:D:997:ARG:HB2	1.95	0.66
1:E:1022:LEU:HD22	1:E:1032:LEU:HD13	1.78	0.65
1:A:1030:PHE:HA	1:A:1036:ARG:O	1.97	0.63
1:A:1036:ARG:HH11	1:A:1036:ARG:HG3	1.62	0.63
1:E:1036:ARG:NH1	1:E:1099:GLN:OE1	2.32	0.63
1:A:998:VAL:HG22	1:A:1123:LEU:HB2	1.81	0.61
1:E:1036:ARG:HD3	1:E:1099:GLN:OE1	2.01	0.61
1:B:1066:LYS:HB2	1:B:1071:LEU:HD11	1.82	0.60
1:D:1036:ARG:NE	1:D:1099:GLN:OE1	2.34	0.60
1:A:1034:LEU:HB3	1:A:1036:ARG:HH11	1.68	0.59
1:A:1122:TYR:HH	1:B:1122:TYR:HH	1.46	0.59
1:F:1010:PRO:HG2	1:F:1104:LEU:HD23	1.85	0.59
1:F:1022:LEU:HD22	1:F:1032:LEU:HD13	1.83	0.58
1:E:1017:ILE:HD12	1:E:1102:LEU:HB3	1.85	0.58
1:F:1041:VAL:HG21	1:F:1126:GLN:HE21	1.70	0.57
1:D:1017:ILE:HD12	1:D:1102:LEU:HB3	1.87	0.56
1:D:997:ARG:HG2	1:D:1123:LEU:HD23	1.87	0.56
1:F:1076:ALA:HB1	1:F:1084:GLU:HG3	1.88	0.56
1:B:1047:PHE:HE2	1:B:1092:LEU:HD11	1.71	0.55
1:A:1105:HIS:CD2	1:A:1106[B]:ARG:HG3	2.42	0.55
1:C:1017:ILE:HD12	1:C:1102:LEU:HB3	1.88	0.54
1:D:996:MET:HB2	1:E:1125:TYR:HB3	1.90	0.54
1:A:1022:LEU:HG	1:A:1032:LEU:HD13	1.90	0.53
1:D:1122:TYR:HH	1:F:1122:TYR:HH	1.52	0.51
1:E:1013:LEU:HD22	1:E:1105:HIS:O	2.11	0.50
1:F:1084:GLU:OE1	4:F:1214:HOH:O	2.20	0.50
1:A:1076:ALA:HB1	1:A:1084:GLU:HB3	1.94	0.49
1:A:998:VAL:HA	1:A:1026:LEU:O	2.13	0.48
1:A:1094:LEU:HD21	1:A:1100:ILE:HD11	1.95	0.47
1:A:1017:ILE:HD12	1:A:1102:LEU:HB3	1.96	0.47
1:D:1122:TYR:OH	1:E:1122:TYR:OH	2.32	0.47
1:B:1047:PHE:CZ	1:B:1100:ILE:HG13	2.50	0.47
1:D:1124:LEU:HD11	1:F:1122:TYR:CD2	2.50	0.46
1:B:1045:TYR:CZ	1:B:1123:LEU:HD13	2.51	0.46
1:A:997:ARG:HG2	1:B:1125:TYR:CD1	2.50	0.46
1:F:1045:TYR:CZ	1:F:1123:LEU:HD13	2.51	0.46
1:E:1122:TYR:HH	1:F:1122:TYR:HH	1.64	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:1071:LEU:HD13	1:F:1092:LEU:HD21	1.98	0.44
1:A:1122:TYR:HH	1:C:1122:TYR:HH	1.52	0.44
1:A:1124:LEU:HD11	1:C:1122:TYR:CD2	2.53	0.43
1:C:1022:LEU:HD13	1:C:1025:ASN:HB2	2.00	0.43
1:A:1031:ASP:OD1	1:A:1034:LEU:HB2	2.18	0.43
1:C:1046:VAL:HG22	1:C:1091:ILE:HG12	2.00	0.43
1:F:1041:VAL:HG21	1:F:1126:GLN:NE2	2.32	0.43
1:E:996:MET:N	4:E:1314:HOH:O	2.51	0.43
1:F:1017:ILE:HD12	1:F:1102:LEU:HB3	2.01	0.43
1:F:1005:THR:HG21	1:F:1021:LEU:HD13	2.01	0.42
1:B:1010:PRO:HG2	1:B:1104:LEU:HD23	2.00	0.42
1:B:1051:MET:SD	1:B:1109:ILE:HD13	2.60	0.42
1:D:1022:LEU:HD22	1:D:1032:LEU:HD13	2.00	0.42
1:D:1010:PRO:HG2	1:D:1104:LEU:HD23	2.01	0.42
1:E:1044:THR:OG1	1:E:1093:GLN:OE1	2.31	0.42
1:B:1076:ALA:HB1	1:B:1084:GLU:HB3	2.00	0.42
1:C:1126:GLN:O	1:C:1127:ASP:HB2	2.19	0.42
1:A:997:ARG:HG2	1:B:1125:TYR:CE1	2.55	0.42
1:F:1019:PHE:HB2	1:F:1032:LEU:HD12	2.02	0.42
1:A:1047:PHE:CZ	1:A:1100:ILE:HG13	2.54	0.42
1:A:1111:GLY:HA2	4:A:1304:HOH:O	2.19	0.42
1:F:1126:GLN:O	1:F:1127:ASP:HB2	2.20	0.42
1:B:1092:LEU:HD13	1:B:1094:LEU:HD21	2.01	0.41
1:C:1053:LYS:HB3	1:C:1060:LEU:HD22	2.01	0.41
1:A:1001:SER:HB3	1:A:1023:LEU:HB2	2.02	0.41
1:D:1036:ARG:HG3	1:D:1101:TRP:HB3	2.02	0.41
1:F:1051:MET:SD	1:F:1109:ILE:HD13	2.60	0.41
1:A:998:VAL:HG22	1:A:1123:LEU:CB	2.50	0.41
1:B:1047:PHE:CE2	1:B:1092:LEU:HD11	2.51	0.41
1:E:1046:VAL:O	1:E:1121:GLY:HA2	2.21	0.41
1:D:997:ARG:HB3	1:D:998:VAL:H	1.69	0.40
1:B:1046:VAL:O	1:B:1121:GLY:HA2	2.21	0.40
1:B:1017:ILE:HD12	1:B:1102:LEU:HB3	2.03	0.40
1:E:1126:GLN:O	1:E:1127:ASP:HB2	2.21	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	131/140 (94%)	122 (93%)	8 (6%)	1 (1%)	16	7
1	B	130/140 (93%)	124 (95%)	6 (5%)	0	100	100
1	C	129/140 (92%)	122 (95%)	7 (5%)	0	100	100
1	D	130/140 (93%)	123 (95%)	6 (5%)	1 (1%)	16	7
1	E	126/140 (90%)	120 (95%)	5 (4%)	1 (1%)	16	7
1	F	130/140 (93%)	122 (94%)	7 (5%)	1 (1%)	16	7
All	All	776/840 (92%)	733 (94%)	39 (5%)	4 (0%)	24	15

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	997	ARG
1	E	997	ARG
1	F	997	ARG
1	A	1030	PHE

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	108/120 (90%)	106 (98%)	2 (2%)	50	40
1	B	102/120 (85%)	99 (97%)	3 (3%)	37	24
1	C	100/120 (83%)	99 (99%)	1 (1%)	68	64

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	104/120 (87%)	100 (96%)	4 (4%)	29	15
1	E	106/120 (88%)	103 (97%)	3 (3%)	38	25
1	F	110/120 (92%)	110 (100%)	0	100	100
All	All	630/720 (88%)	617 (98%)	13 (2%)	47	37

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	998	VAL
1	A	1084	GLU
1	B	1028	GLU
1	B	1084	GLU
1	B	1092	LEU
1	C	1058	VAL
1	D	1022	LEU
1	D	1068	GLU
1	D	1084	GLU
1	D	1106	ARG
1	E	997	ARG
1	E	1012	THR
1	E	1022	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1015	GLN
1	A	1038	ASN
1	B	1024	ASN
1	B	1126	GLN
1	E	1077	ASN
1	F	1033	GLN
1	F	1038	ASN

5.3.3 RNA ⓘ

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 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 for Mogul analysis.

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	GOL	D	1202	-	5,5,5	0.36	0	5,5,5	0.73	0

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	GOL	D	1202	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	1202	GOL	O1-C1-C2-C3
3	D	1202	GOL	C1-C2-C3-O3
3	D	1202	GOL	O1-C1-C2-O2
3	D	1202	GOL	O2-C2-C3-O3

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

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

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	132/140 (94%)	0.56	11 (8%) 17 19	17, 35, 51, 60	1 (0%)
1	B	132/140 (94%)	0.55	6 (4%) 38 43	21, 36, 50, 58	0
1	C	131/140 (93%)	0.57	4 (3%) 51 57	20, 36, 52, 59	0
1	D	132/140 (94%)	0.67	8 (6%) 27 30	18, 35, 52, 56	0
1	E	130/140 (92%)	0.59	5 (3%) 44 49	17, 36, 53, 58	0
1	F	132/140 (94%)	0.65	10 (7%) 20 22	19, 37, 55, 60	0
All	All	789/840 (93%)	0.60	44 (5%) 30 33	17, 36, 53, 60	1 (0%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	996	MET	4.7
1	A	1128	LEU	4.1
1	D	1032	LEU	3.5
1	D	996	MET	3.4
1	A	1127	ASP	3.3
1	F	996	MET	3.2
1	C	1031	ASP	3.1
1	E	1079	GLY	2.9
1	B	1022	LEU	2.9
1	C	1032	LEU	2.8
1	F	1032	LEU	2.7
1	A	1011	GLY	2.7
1	F	1010	PRO	2.6
1	D	1079	GLY	2.6
1	B	1124	LEU	2.6
1	A	1010	PRO	2.5
1	F	1008	LEU	2.5
1	B	1032	LEU	2.4
1	F	1011	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	F	1043	GLY	2.4
1	A	1025	ASN	2.3
1	A	1031	ASP	2.3
1	C	1041	VAL	2.3
1	C	1034	LEU	2.3
1	F	1012	THR	2.3
1	D	1127	ASP	2.3
1	B	1080	ALA	2.2
1	D	1106	ARG	2.2
1	B	1009	ALA	2.2
1	F	1080	ALA	2.2
1	E	1013	LEU	2.2
1	E	1095	PHE	2.1
1	A	1096	GLN	2.1
1	A	1036	ARG	2.1
1	E	1032	LEU	2.1
1	D	1057	ASN	2.1
1	A	1030	PHE	2.1
1	A	1032	LEU	2.1
1	D	1010	PRO	2.1
1	F	1039	CYS	2.1
1	D	1011	GLY	2.0
1	A	1124	LEU	2.0
1	B	1008	LEU	2.0
1	F	1079	GLY	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no oligosaccharides in this entry.

6.4 Ligands

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. 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	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	D	1202	6/6	0.78	0.25	20,20,20,20	0
2	CA	E	1201	1/1	0.96	0.12	55,55,55,55	0
2	CA	B	1201	1/1	0.98	0.10	38,38,38,38	0
2	CA	A	1201	1/1	0.99	0.02	24,24,24,24	0
2	CA	D	1201	1/1	0.99	0.03	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.