



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:01 AM GMT

PDB ID : 2F6A
Title : Collagen Adhesin and Collagen Complex Structure
Authors : Zong, Y.; Narayana, S.L.V.
Deposited on : 2005-11-28
Resolution : 3.29 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

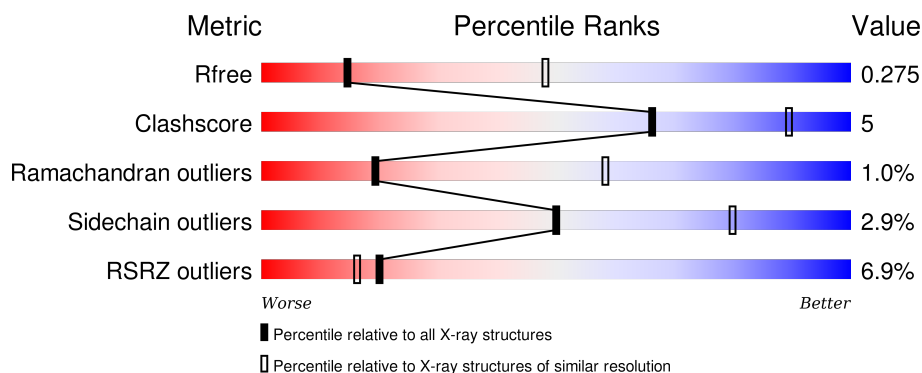
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.29 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}	91344	2060 (3.40-3.20)
Clashscore	102246	1058 (3.38-3.22)
Ramachandran outliers	100387	1038 (3.38-3.22)
Sidechain outliers	100360	1037 (3.38-3.22)
RSRZ outliers	91569	2070 (3.40-3.20)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	303	<div> <div>3%</div> <div>84%</div> <div>15%</div> <div>.</div> </div>
1	B	303	<div> <div>12%</div> <div>86%</div> <div>13%</div> <div>.</div> </div>
1	C	303	<div> <div>5%</div> <div>85%</div> <div>13%</div> <div>..</div> </div>
1	D	303	<div> <div>6%</div> <div>84%</div> <div>13%</div> <div>..</div> </div>
2	E	30	<div> <div>10%</div> <div>97%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
2	F	30	
2	G	30	
2	H	30	
2	I	30	
2	J	30	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	HYP	E	12	X	-	-	-
2	HYP	E	21	X	-	-	-
2	HYP	E	24	X	-	-	-
2	HYP	E	27	X	-	-	-
2	HYP	E	3	X	-	-	-
2	HYP	E	30	X	-	-	-
2	HYP	E	6	X	-	-	-
2	HYP	E	9	X	-	-	-
2	HYP	F	12	X	-	-	-
2	HYP	F	21	X	-	-	-
2	HYP	F	24	X	-	-	-
2	HYP	F	27	X	-	-	-
2	HYP	F	3	X	-	-	-
2	HYP	F	30	X	-	-	-
2	HYP	F	6	X	-	-	-
2	HYP	F	9	X	-	-	-
2	HYP	G	12	X	-	-	-
2	HYP	G	21	X	-	-	-
2	HYP	G	24	X	-	-	-
2	HYP	G	27	X	-	-	-
2	HYP	G	3	X	-	-	-
2	HYP	G	30	X	-	-	-
2	HYP	G	6	X	-	-	-
2	HYP	G	9	X	-	-	-
2	HYP	H	12	X	-	-	-
2	HYP	H	21	X	-	-	-
2	HYP	H	24	X	-	-	-
2	HYP	H	27	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	HYP	H	3	X	-	-	-
2	HYP	H	30	X	-	-	-
2	HYP	H	6	X	-	-	-
2	HYP	H	9	X	-	-	-
2	HYP	I	12	X	-	-	-
2	HYP	I	21	X	-	-	-
2	HYP	I	24	X	-	-	-
2	HYP	I	27	X	-	-	-
2	HYP	I	3	X	-	-	-
2	HYP	I	30	X	-	-	-
2	HYP	I	6	X	-	-	-
2	HYP	I	9	X	-	-	-
2	HYP	J	12	X	-	-	-
2	HYP	J	21	X	-	-	-
2	HYP	J	24	X	-	-	-
2	HYP	J	27	X	-	-	-
2	HYP	J	3	X	-	-	-
2	HYP	J	6	X	-	-	-
2	HYP	J	9	X	-	-	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10394 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 Collagen adhesin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	303	Total	C	N	O	S	0	0	0
			2327	1442	394	488	3			
1	B	299	Total	C	N	O	S	0	0	0
			2302	1428	389	482	3			
1	C	299	Total	C	N	O	S	0	0	0
			2302	1428	389	482	3			
1	D	299	Total	C	N	O	S	0	0	0
			2302	1428	389	482	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	28	GLY	-	CLONING ARTIFACT	UNP Q53654
A	29	SER	-	CLONING ARTIFACT	UNP Q53654
B	28	GLY	-	CLONING ARTIFACT	UNP Q53654
B	29	SER	-	CLONING ARTIFACT	UNP Q53654
C	28	GLY	-	CLONING ARTIFACT	UNP Q53654
C	29	SER	-	CLONING ARTIFACT	UNP Q53654
D	28	GLY	-	CLONING ARTIFACT	UNP Q53654
D	29	SER	-	CLONING ARTIFACT	UNP Q53654

- Molecule 2 is a protein called Collagen.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	30	Total	C	N	O	0	0	0
			196	121	36	39			
2	F	30	Total	C	N	O	0	0	0
			196	121	36	39			
2	G	30	Total	C	N	O	0	0	0
			196	121	36	39			
2	H	30	Total	C	N	O	0	0	0
			196	121	36	39			

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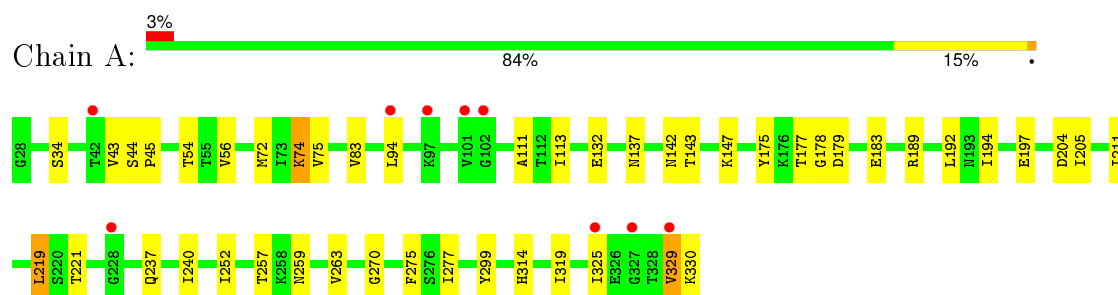
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	I	30	Total	C	N	O	0	0	0
			196	121	36	39			
2	J	28	Total	C	N	O	0	0	0
			181	111	34	36			

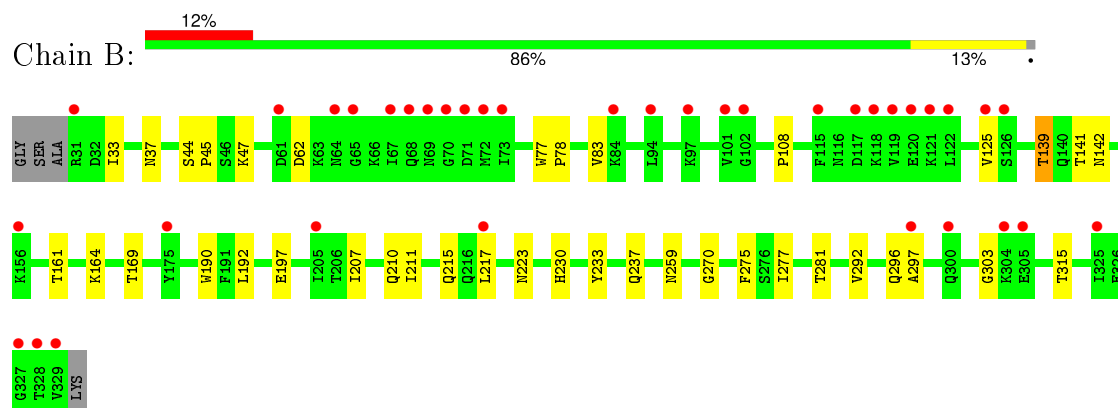
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

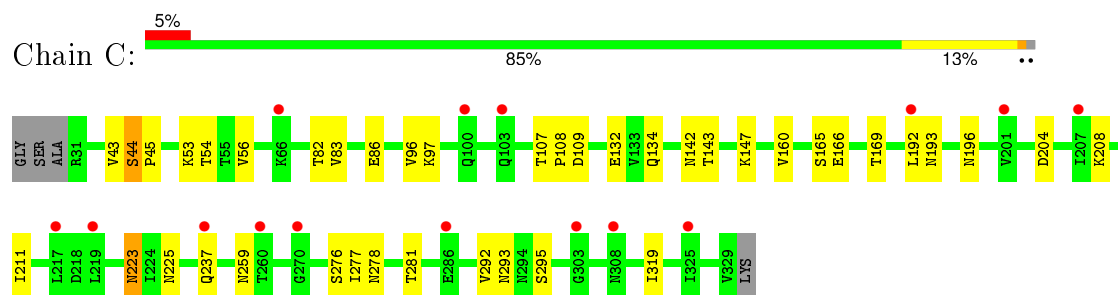
- Molecule 1: Collagen adhesin



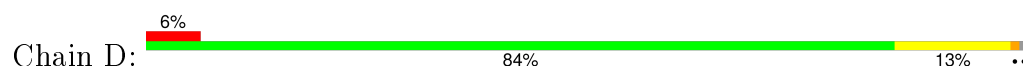
- Molecule 1: Collagen adhesin

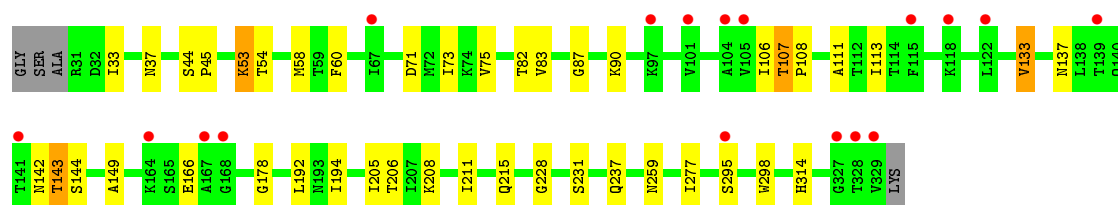


- Molecule 1: Collagen adhesin

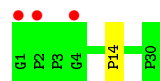


- Molecule 1: Collagen adhesin

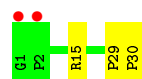
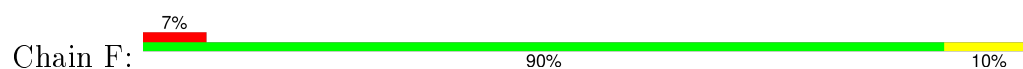




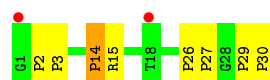
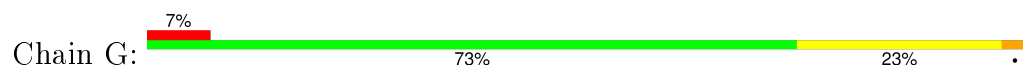
- Molecule 2: Collagen



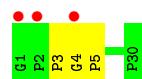
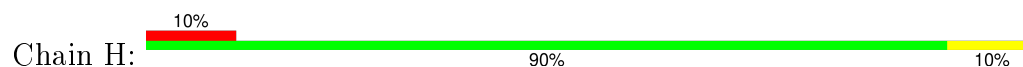
- Molecule 2: Collagen



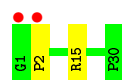
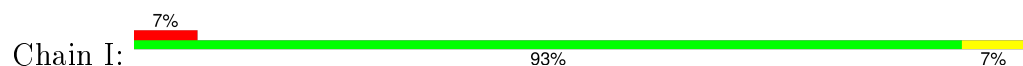
- Molecule 2: Collagen



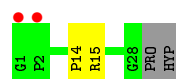
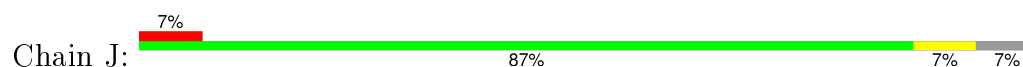
- Molecule 2: Collagen



- Molecule 2: Collagen



- Molecule 2: Collagen



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	90.55Å 193.82Å 205.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.29 46.80 – 3.29	Depositor EDS
% Data completeness (in resolution range)	94.5 (50.00-3.29) 94.4 (46.80-3.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 3.32Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.252 , 0.308 0.274 , 0.275	Depositor DCC
R_{free} test set	1308 reflections (4.99%)	DCC
Wilson B-factor (Å ²)	77.4	Xtriage
Anisotropy	0.695	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 44.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 26252 reflections	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	10394	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: HYP

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	0.31	0/2366	0.44	0/3205
1	B	0.32	0/2341	0.44	0/3174
1	C	0.32	0/2341	0.44	0/3174
1	D	0.32	0/2341	0.44	0/3174
2	E	0.36	0/133	0.59	0/172
2	F	0.35	0/133	0.54	0/172
2	G	0.37	0/133	0.54	0/172
2	H	0.35	0/133	0.55	0/172
2	I	0.36	0/133	0.52	0/172
2	J	0.35	0/125	0.53	0/160
All	All	0.32	0/10179	0.45	0/13747

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
2	E	8	0
2	F	8	0
2	G	8	0
2	H	8	0
2	I	8	0
2	J	7	0
All	All	47	0

There are no bond length outliers.

There are no bond angle outliers.

All (47) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	E	3	HYP	CG
2	E	6	HYP	CG
2	E	9	HYP	CG
2	E	12	HYP	CG
2	E	21	HYP	CG
2	E	24	HYP	CG
2	E	27	HYP	CG
2	E	30	HYP	CG
2	F	3	HYP	CG
2	F	6	HYP	CG
2	F	9	HYP	CG
2	F	12	HYP	CG
2	F	21	HYP	CG
2	F	24	HYP	CG
2	F	27	HYP	CG
2	F	30	HYP	CG
2	G	3	HYP	CG
2	G	6	HYP	CG
2	G	9	HYP	CG
2	G	12	HYP	CG
2	G	21	HYP	CG
2	G	24	HYP	CG
2	G	27	HYP	CG
2	G	30	HYP	CG
2	H	3	HYP	CG
2	H	6	HYP	CG
2	H	9	HYP	CG
2	H	12	HYP	CG
2	H	21	HYP	CG
2	H	24	HYP	CG
2	H	27	HYP	CG
2	H	30	HYP	CG
2	I	3	HYP	CG
2	I	6	HYP	CG
2	I	9	HYP	CG
2	I	12	HYP	CG
2	I	21	HYP	CG
2	I	24	HYP	CG
2	I	27	HYP	CG
2	I	30	HYP	CG
2	J	3	HYP	CG
2	J	6	HYP	CG
2	J	9	HYP	CG

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Mol	Chain	Res	Type	Atom
2	J	12	HYP	CG
2	J	21	HYP	CG
2	J	24	HYP	CG
2	J	27	HYP	CG

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	2327	0	2254	24	0
1	B	2302	0	2228	21	0
1	C	2302	0	2228	20	0
1	D	2302	0	2228	25	0
2	E	196	0	184	1	0
2	F	196	0	184	2	0
2	G	196	0	184	7	0
2	H	196	0	184	2	0
2	I	196	0	184	2	0
2	J	181	0	170	1	0
All	All	10394	0	10028	99	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 (99) 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:44:SER:HB3	1:C:45:PRO:HD3	1.46	0.94
1:A:44:SER:HB3	1:A:45:PRO:HD3	1.58	0.85
1:D:58:MET:HE3	1:D:73:ILE:HG21	1.68	0.75
1:B:44:SER:HB3	1:B:45:PRO:HD3	1.73	0.71
1:D:211:ILE:HB	1:D:259:ASN:HB3	1.74	0.69
1:B:190:TRP:HE1	1:B:215:GLN:HE22	1.40	0.69
1:C:223:ASN:HD21	1:C:278:ASN:HB3	1.59	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:228:GLY:H	1:D:231:SER:HB3	1.62	0.62
1:C:45:PRO:HD2	1:C:54:THR:HG23	1.80	0.62
1:B:141:THR:HG22	1:B:142:ASN:N	2.14	0.62
1:C:165:SER:HB3	2:G:29:PRO:HG3	1.80	0.62
1:D:192:LEU:HB2	1:D:277:ILE:HB	1.82	0.61
1:C:44:SER:CB	1:C:45:PRO:HD3	2.25	0.60
1:B:210:GLN:HE22	1:B:296:GLN:HE22	1.51	0.58
1:D:194:ILE:HG21	1:D:205:ILE:HG21	1.87	0.56
1:C:225:ASN:HB3	1:C:276:SER:HB3	1.88	0.56
2:G:29:PRO:HB2	2:G:30:HYP:HD22	1.88	0.56
1:D:54:THR:HB	1:D:133:VAL:HG12	1.87	0.55
1:B:211:ILE:HB	1:B:259:ASN:HB3	1.89	0.55
1:C:211:ILE:HB	1:C:259:ASN:HB3	1.87	0.55
1:C:43:VAL:HG11	1:C:160:VAL:HG22	1.89	0.55
1:C:192:LEU:HB2	1:C:277:ILE:HB	1.89	0.54
1:C:82:THR:HG21	1:C:147:LYS:HE2	1.89	0.54
1:B:192:LEU:HB2	1:B:277:ILE:HB	1.89	0.54
1:D:33:ILE:HB	1:D:37:ASN:HD22	1.73	0.54
1:D:45:PRO:HG2	1:D:54:THR:HA	1.90	0.53
1:D:58:MET:HE2	1:D:60:PHE:HB3	1.90	0.53
1:A:192:LEU:HB2	1:A:277:ILE:HB	1.92	0.52
1:A:132:GLU:HB2	1:A:319:ILE:HG23	1.91	0.51
1:D:178:GLY:H	1:D:314:HIS:CD2	2.28	0.51
1:C:44:SER:HB3	1:C:45:PRO:CD	2.29	0.50
1:A:94:LEU:HD22	1:A:325:ILE:HB	1.93	0.50
1:A:329:VAL:HB	1:A:330:LYS:HD3	1.93	0.49
1:C:193:ASN:HB3	1:C:196:ASN:OD1	2.13	0.49
1:A:179:ASP:HB3	1:A:189:ARG:HE	1.77	0.49
1:B:190:TRP:HE1	1:B:215:GLN:NE2	2.08	0.49
1:A:44:SER:CB	1:A:45:PRO:HD3	2.38	0.49
1:A:194:ILE:HB	1:A:275:PHE:HB2	1.95	0.49
1:A:34:SER:HB3	1:A:74:LYS:H	1.77	0.48
1:A:270:GLY:HA2	1:A:275:PHE:HE1	1.78	0.48
1:D:82:THR:HG23	1:D:83:VAL:HG23	1.96	0.48
1:B:217:LEU:HA	1:B:281:THR:HG22	1.96	0.48
1:A:75:VAL:HB	1:A:111:ALA:HB3	1.95	0.48
1:B:141:THR:HG22	1:B:142:ASN:H	1.75	0.48
1:C:86:GLU:HB2	1:C:134:GLN:HB3	1.95	0.47
2:I:15:ARG:HG2	2:J:14:PRO:HG2	1.95	0.47
1:B:77:TRP:HB2	1:B:78:PRO:HD2	1.97	0.47
1:D:87:GLY:HA2	1:D:133:VAL:HG23	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:139:THR:O	1:B:164:LYS:HG2	2.14	0.47
1:C:132:GLU:HB2	1:C:319:ILE:HG23	1.97	0.46
2:G:29:PRO:CB	2:G:30:HYP:HD22	2.45	0.46
1:A:175:TYR:HE2	1:A:177:THR:HG1	1.63	0.46
1:C:43:VAL:HG22	1:C:56:VAL:HG22	1.98	0.46
1:D:44:SER:OG	1:D:45:PRO:HD3	2.15	0.46
1:B:33:ILE:HB	1:B:37:ASN:HD22	1.79	0.45
1:D:107:THR:HG22	1:D:108:PRO:HD2	1.98	0.45
1:A:240:ILE:HG23	1:A:252:ILE:HB	1.98	0.45
1:A:211:ILE:HB	1:A:259:ASN:HB3	1.99	0.45
1:D:33:ILE:HD11	1:D:71:ASP:HB3	1.99	0.45
1:D:143:THR:HG22	1:D:166:GLU:HG2	1.99	0.45
1:B:270:GLY:HA2	1:B:275:PHE:HE1	1.81	0.44
2:F:15:ARG:HG3	2:G:14:PRO:HB2	1.98	0.44
1:A:219:LEU:HD13	1:B:230:HIS:CE1	2.52	0.44
1:D:143:THR:HB	1:D:144:SER:H	1.64	0.44
1:A:205:ILE:HG23	1:A:299:TYR:HB3	1.99	0.44
2:H:3:HYP:HA	2:I:2:PRO:HB2	2.00	0.44
1:B:47:LYS:HG2	1:B:161:THR:HB	2.00	0.43
1:A:178:GLY:H	1:A:314:HIS:CD2	2.36	0.43
1:B:141:THR:CG2	1:B:142:ASN:N	2.80	0.43
1:A:270:GLY:HA2	1:A:275:PHE:CE1	2.53	0.43
1:A:45:PRO:HD2	1:A:54:THR:HG23	2.00	0.43
1:A:43:VAL:HG22	1:A:56:VAL:HG22	2.01	0.43
1:C:44:SER:CB	1:C:45:PRO:CD	2.95	0.43
1:B:292:VAL:HG22	1:B:315:THR:HG23	2.01	0.43
1:C:223:ASN:HD22	1:C:223:ASN:C	2.22	0.43
2:G:2:PRO:HA	2:G:3:HYP:HD23	1.94	0.42
1:A:44:SER:HB3	1:A:45:PRO:CD	2.40	0.42
1:D:206:THR:HB	1:D:298:TRP:HB2	2.00	0.42
1:A:72:MET:HA	1:A:113:ILE:O	2.19	0.42
1:D:211:ILE:HG23	1:D:215:GLN:HE21	1.85	0.42
2:G:26:PRO:HA	2:G:27:HYP:HD23	1.96	0.42
1:C:107:THR:C	1:C:109:ASP:H	2.23	0.42
1:D:75:VAL:HB	1:D:111:ALA:HB3	2.01	0.42
1:D:90:LYS:HB3	1:D:106:ILE:HD12	2.02	0.42
1:D:83:VAL:HG11	1:D:149:ALA:HB2	2.01	0.41
1:A:252:ILE:HG23	1:A:263:VAL:HG22	2.02	0.41
1:B:223:ASN:HD22	1:B:233:TYR:HE1	1.67	0.41
1:D:53:LYS:H	1:D:53:LYS:HD3	1.85	0.41
2:F:29:PRO:HA	2:F:30:HYP:HD23	1.97	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:96:VAL:HG22	1:C:97:LYS:HG2	2.02	0.41
1:C:208:LYS:O	1:C:295:SER:HA	2.21	0.41
1:B:207:ILE:HG22	1:B:297:ALA:HA	2.02	0.41
1:A:175:TYR:CE1	2:E:14:PRO:HG3	2.56	0.41
1:D:208:LYS:O	1:D:295:SER:HA	2.21	0.41
2:G:29:PRO:HB2	2:G:30:HYP:CD	2.50	0.40
1:B:33:ILE:HB	1:B:37:ASN:ND2	2.36	0.40
2:H:4:GLY:HA2	2:H:5:PRO:HD3	1.91	0.40
1:D:73:ILE:HB	1:D:113:ILE:HD12	2.03	0.40
1:B:62:ASP:HB3	1:B:125:VAL:HB	2.03	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	301/303 (99%)	279 (93%)	21 (7%)	1 (0%)	46	81
1	B	297/303 (98%)	277 (93%)	16 (5%)	4 (1%)	15	52
1	C	297/303 (98%)	276 (93%)	16 (5%)	5 (2%)	11	47
1	D	297/303 (98%)	275 (93%)	20 (7%)	2 (1%)	26	66
2	E	21/30 (70%)	21 (100%)	0	0	100	100
2	F	21/30 (70%)	21 (100%)	0	0	100	100
2	G	21/30 (70%)	19 (90%)	1 (5%)	1 (5%)	3	20
2	H	21/30 (70%)	21 (100%)	0	0	100	100
2	I	21/30 (70%)	21 (100%)	0	0	100	100
2	J	19/30 (63%)	19 (100%)	0	0	100	100
All	All	1316/1392 (94%)	1229 (93%)	74 (6%)	13 (1%)	19	58

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	142	ASN
1	D	142	ASN
1	A	237	GLN
1	B	139	THR
2	G	14	PRO
1	B	237	GLN
1	B	303	GLY
1	C	237	GLN
1	C	108	PRO
1	C	44	SER
1	C	166	GLU
1	D	237	GLN
1	B	108	PRO

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	264/265 (100%)	251 (95%)	13 (5%)	31	69
1	B	262/265 (99%)	259 (99%)	3 (1%)	80	90
1	C	262/265 (99%)	253 (97%)	9 (3%)	44	77
1	D	262/265 (99%)	257 (98%)	5 (2%)	65	85
2	E	12/12 (100%)	12 (100%)	0	100	100
2	F	12/12 (100%)	12 (100%)	0	100	100
2	G	12/12 (100%)	11 (92%)	1 (8%)	14	47
2	H	12/12 (100%)	12 (100%)	0	100	100
2	I	12/12 (100%)	12 (100%)	0	100	100
2	J	11/12 (92%)	10 (91%)	1 (9%)	12	41
All	All	1121/1132 (99%)	1089 (97%)	32 (3%)	50	80

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	LYS
1	A	83	VAL
1	A	137	ASN
1	A	142	ASN
1	A	143	THR
1	A	147	LYS
1	A	183	GLU
1	A	197	GLU
1	A	204	ASP
1	A	219	LEU
1	A	221	THR
1	A	257	THR
1	A	329	VAL
1	B	83	VAL
1	B	169	THR
1	B	197	GLU
1	C	53	LYS
1	C	83	VAL
1	C	143	THR
1	C	169	THR
1	C	204	ASP
1	C	223	ASN
1	C	281	THR
1	C	292	VAL
1	C	293	ASN
1	D	53	LYS
1	D	107	THR
1	D	133	VAL
1	D	137	ASN
1	D	143	THR
2	G	15	ARG
2	J	15	ARG

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

Mol	Chain	Res	Type
1	A	37	ASN
1	A	137	ASN
1	A	155	ASN
1	A	210	GLN
1	A	215	GLN
1	A	223	ASN
1	A	278	ASN

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Mol	Chain	Res	Type
1	A	296	GLN
1	A	314	HIS
1	B	37	ASN
1	B	68	GLN
1	B	69	ASN
1	B	137	ASN
1	B	155	ASN
1	B	210	GLN
1	B	215	GLN
1	B	223	ASN
1	B	230	HIS
1	C	37	ASN
1	C	69	ASN
1	C	155	ASN
1	C	215	GLN
1	C	223	ASN
1	C	293	ASN
1	C	294	ASN
1	C	314	HIS
1	D	37	ASN
1	D	137	ASN
1	D	155	ASN
1	D	278	ASN
1	D	293	ASN
1	D	313	ASN
1	D	314	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

47 non-standard protein/DNA/RNA residues are modelled in this entry.

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
2	HYP	E	12	2	7,8,9	0.54	0	5,10,12	1.88	2 (40%)
2	HYP	E	21	2	7,8,9	0.55	0	5,10,12	2.19	2 (40%)
2	HYP	E	24	2	7,8,9	0.55	0	5,10,12	1.64	1 (20%)
2	HYP	E	27	2	7,8,9	0.57	0	5,10,12	1.50	1 (20%)
2	HYP	E	3	2	7,8,9	0.55	0	5,10,12	2.06	2 (40%)
2	HYP	E	30	2	7,8,9	0.52	0	5,10,12	2.03	3 (60%)
2	HYP	E	6	2	7,8,9	0.54	0	5,10,12	2.20	2 (40%)
2	HYP	E	9	2	7,8,9	0.56	0	5,10,12	1.60	1 (20%)
2	HYP	F	12	2	7,8,9	0.56	0	5,10,12	2.05	2 (40%)
2	HYP	F	21	2	7,8,9	0.55	0	5,10,12	2.11	2 (40%)
2	HYP	F	24	2	7,8,9	0.55	0	5,10,12	2.15	2 (40%)
2	HYP	F	27	2	7,8,9	0.55	0	5,10,12	1.76	2 (40%)
2	HYP	F	3	2	7,8,9	0.57	0	5,10,12	2.15	2 (40%)
2	HYP	F	30	2	7,8,9	0.54	0	5,10,12	2.18	2 (40%)
2	HYP	F	6	2	7,8,9	0.55	0	5,10,12	2.15	2 (40%)
2	HYP	F	9	2	7,8,9	0.55	0	5,10,12	2.08	2 (40%)
2	HYP	G	12	2	7,8,9	0.57	0	5,10,12	2.18	2 (40%)
2	HYP	G	21	2	7,8,9	0.59	0	5,10,12	2.00	3 (60%)
2	HYP	G	24	2	7,8,9	0.56	0	5,10,12	2.17	2 (40%)
2	HYP	G	27	2	7,8,9	0.55	0	5,10,12	1.77	2 (40%)
2	HYP	G	3	2	7,8,9	0.54	0	5,10,12	2.10	2 (40%)
2	HYP	G	30	2	7,8,9	0.55	0	5,10,12	1.56	2 (40%)
2	HYP	G	6	2	7,8,9	0.57	0	5,10,12	2.07	3 (60%)
2	HYP	G	9	2	7,8,9	0.57	0	5,10,12	2.13	3 (60%)
2	HYP	H	12	2	7,8,9	0.55	0	5,10,12	1.53	0
2	HYP	H	21	2	7,8,9	0.55	0	5,10,12	2.00	2 (40%)
2	HYP	H	24	2	7,8,9	0.52	0	5,10,12	1.75	2 (40%)
2	HYP	H	27	2	7,8,9	0.59	0	5,10,12	1.89	2 (40%)
2	HYP	H	3	2	7,8,9	0.54	0	5,10,12	2.10	2 (40%)
2	HYP	H	30	2	7,8,9	0.52	0	5,10,12	2.15	3 (60%)
2	HYP	H	6	2	7,8,9	0.53	0	5,10,12	2.00	2 (40%)
2	HYP	H	9	2	7,8,9	0.54	0	5,10,12	1.76	2 (40%)
2	HYP	I	12	2	7,8,9	0.56	0	5,10,12	2.11	2 (40%)
2	HYP	I	21	2	7,8,9	0.56	0	5,10,12	1.97	3 (60%)
2	HYP	I	24	2	7,8,9	0.54	0	5,10,12	2.14	2 (40%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HYP	I	27	2	7,8,9	0.56	0	5,10,12	1.72	2 (40%)
2	HYP	I	3	2	7,8,9	0.58	0	5,10,12	2.13	2 (40%)
2	HYP	I	30	2	7,8,9	0.53	0	5,10,12	2.22	2 (40%)
2	HYP	I	6	2	7,8,9	0.55	0	5,10,12	2.07	2 (40%)
2	HYP	I	9	2	7,8,9	0.56	0	5,10,12	2.15	2 (40%)
2	HYP	J	12	2	7,8,9	0.54	0	5,10,12	2.08	2 (40%)
2	HYP	J	21	2	7,8,9	0.59	0	5,10,12	2.05	3 (60%)
2	HYP	J	24	2	7,8,9	0.56	0	5,10,12	2.17	3 (60%)
2	HYP	J	27	2	7,8,9	0.57	0	5,10,12	2.12	2 (40%)
2	HYP	J	3	2	7,8,9	0.55	0	5,10,12	2.07	2 (40%)
2	HYP	J	6	2	7,8,9	0.60	0	5,10,12	2.17	3 (60%)
2	HYP	J	9	2	7,8,9	0.56	0	5,10,12	2.18	2 (40%)

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
2	HYP	E	12	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	E	21	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	E	24	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	E	27	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	E	3	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	E	30	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	E	6	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	E	9	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	F	12	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	F	21	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	F	24	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	F	27	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	F	3	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	F	30	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	F	6	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	F	9	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	G	12	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	G	21	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	G	24	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	G	27	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	G	3	2	1/1/2/4	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HYP	G	30	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	G	6	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	G	9	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	H	12	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	H	21	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	H	24	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	H	27	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	H	3	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	H	30	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	H	6	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	H	9	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	I	12	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	I	21	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	I	24	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	I	27	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	I	3	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	I	30	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	I	6	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	I	9	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	J	12	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	J	21	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	J	24	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	J	27	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	J	3	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	J	6	2	1/1/2/4	0/0/11/13	0/1/1/1
2	HYP	J	9	2	1/1/2/4	0/0/11/13	0/1/1/1

There are no bond length outliers.

All (98) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	24	HYP	O-C-CA	-2.34	119.27	125.44
2	E	30	HYP	O-C-CA	-2.09	119.92	125.44
2	H	30	HYP	O-C-CA	-2.08	119.94	125.44
2	J	6	HYP	O-C-CA	-2.08	119.95	125.44
2	G	30	HYP	O-C-CA	-2.08	119.95	125.44
2	G	21	HYP	O-C-CA	-2.06	120.01	125.44
2	J	21	HYP	O-C-CA	-2.05	120.03	125.44
2	I	21	HYP	O-C-CA	-2.03	120.08	125.44
2	G	6	HYP	O-C-CA	-2.02	120.10	125.44
2	J	24	HYP	O-C-CA	-2.01	120.14	125.44
2	G	9	HYP	O-C-CA	-2.00	120.16	125.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	27	HYP	CG-CB-CA	2.05	106.83	103.90
2	E	27	HYP	CB-CG-CD	2.10	105.73	103.14
2	F	27	HYP	CG-CB-CA	2.19	107.04	103.90
2	E	12	HYP	CG-CB-CA	2.21	107.06	103.90
2	E	9	HYP	CB-CG-CD	2.25	105.92	103.14
2	H	27	HYP	CG-CB-CA	2.26	107.14	103.90
2	H	9	HYP	CG-CB-CA	2.30	107.19	103.90
2	G	21	HYP	CG-CB-CA	2.31	107.21	103.90
2	H	21	HYP	CG-CB-CA	2.32	107.23	103.90
2	J	27	HYP	CG-CB-CA	2.35	107.27	103.90
2	J	12	HYP	CG-CB-CA	2.36	107.28	103.90
2	G	27	HYP	CG-CB-CA	2.37	107.29	103.90
2	J	21	HYP	CG-CB-CA	2.47	107.44	103.90
2	G	27	HYP	CB-CG-CD	2.49	106.21	103.14
2	G	30	HYP	CG-CB-CA	2.49	107.46	103.90
2	H	9	HYP	CB-CG-CD	2.49	106.22	103.14
2	G	12	HYP	CG-CB-CA	2.51	107.50	103.90
2	I	21	HYP	CG-CB-CA	2.53	107.52	103.90
2	E	24	HYP	CB-CG-CD	2.58	106.33	103.14
2	I	27	HYP	CB-CG-CD	2.58	106.33	103.14
2	J	6	HYP	CG-CB-CA	2.58	107.60	103.90
2	F	27	HYP	CB-CG-CD	2.61	106.37	103.14
2	I	12	HYP	CG-CB-CA	2.62	107.66	103.90
2	F	9	HYP	CG-CB-CA	2.68	107.73	103.90
2	H	6	HYP	CG-CB-CA	2.68	107.74	103.90
2	G	6	HYP	CG-CB-CA	2.69	107.75	103.90
2	J	3	HYP	CG-CB-CA	2.69	107.76	103.90
2	E	3	HYP	CG-CB-CA	2.71	107.78	103.90
2	F	3	HYP	CG-CB-CA	2.72	107.80	103.90
2	I	24	HYP	CG-CB-CA	2.79	107.89	103.90
2	F	12	HYP	CG-CB-CA	2.79	107.90	103.90
2	E	30	HYP	CB-CG-CD	2.79	106.59	103.14
2	I	6	HYP	CG-CB-CA	2.80	107.91	103.90
2	F	21	HYP	CG-CB-CA	2.81	107.92	103.90
2	G	3	HYP	CG-CB-CA	2.82	107.93	103.90
2	H	3	HYP	CG-CB-CA	2.82	107.94	103.90
2	I	9	HYP	CG-CB-CA	2.83	107.95	103.90
2	E	21	HYP	CG-CB-CA	2.85	107.99	103.90
2	E	30	HYP	CG-CB-CA	2.86	107.99	103.90
2	G	9	HYP	CG-CB-CA	2.87	108.02	103.90
2	G	24	HYP	CG-CB-CA	2.89	108.03	103.90
2	J	9	HYP	CG-CB-CA	2.89	108.04	103.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	24	HYP	CG-CB-CA	2.90	108.05	103.90
2	I	3	HYP	CB-CG-CD	2.90	106.72	103.14
2	F	6	HYP	CG-CB-CA	2.90	108.06	103.90
2	J	24	HYP	CG-CB-CA	2.92	108.08	103.90
2	H	27	HYP	CB-CG-CD	2.92	106.75	103.14
2	F	30	HYP	CG-CB-CA	2.94	108.11	103.90
2	I	21	HYP	CB-CG-CD	2.96	106.80	103.14
2	H	6	HYP	CB-CG-CD	2.97	106.81	103.14
2	I	30	HYP	CG-CB-CA	2.98	108.17	103.90
2	E	12	HYP	CB-CG-CD	2.98	106.83	103.14
2	E	6	HYP	CG-CB-CA	3.03	108.23	103.90
2	H	30	HYP	CG-CB-CA	3.04	108.25	103.90
2	F	12	HYP	CB-CG-CD	3.04	106.89	103.14
2	H	24	HYP	CB-CG-CD	3.05	106.91	103.14
2	H	30	HYP	CB-CG-CD	3.05	106.92	103.14
2	I	6	HYP	CB-CG-CD	3.11	106.98	103.14
2	E	3	HYP	CB-CG-CD	3.16	107.05	103.14
2	G	3	HYP	CB-CG-CD	3.16	107.05	103.14
2	H	3	HYP	CB-CG-CD	3.17	107.06	103.14
2	J	3	HYP	CB-CG-CD	3.18	107.07	103.14
2	G	6	HYP	CB-CG-CD	3.18	107.07	103.14
2	G	21	HYP	CB-CG-CD	3.20	107.10	103.14
2	G	9	HYP	CB-CG-CD	3.21	107.11	103.14
2	F	21	HYP	CB-CG-CD	3.22	107.12	103.14
2	F	9	HYP	CB-CG-CD	3.22	107.13	103.14
2	H	21	HYP	CB-CG-CD	3.23	107.14	103.14
2	I	3	HYP	CG-CB-CA	3.25	108.55	103.90
2	J	21	HYP	CB-CG-CD	3.26	107.16	103.14
2	J	24	HYP	CB-CG-CD	3.27	107.18	103.14
2	F	6	HYP	CB-CG-CD	3.27	107.18	103.14
2	F	24	HYP	CB-CG-CD	3.30	107.22	103.14
2	E	6	HYP	CB-CG-CD	3.31	107.23	103.14
2	G	24	HYP	CB-CG-CD	3.33	107.25	103.14
2	I	9	HYP	CB-CG-CD	3.33	107.26	103.14
2	F	30	HYP	CB-CG-CD	3.33	107.26	103.14
2	I	24	HYP	CB-CG-CD	3.35	107.28	103.14
2	I	12	HYP	CB-CG-CD	3.36	107.29	103.14
2	J	9	HYP	CB-CG-CD	3.36	107.30	103.14
2	F	3	HYP	CB-CG-CD	3.40	107.35	103.14
2	I	30	HYP	CB-CG-CD	3.41	107.36	103.14
2	E	21	HYP	CB-CG-CD	3.44	107.39	103.14
2	J	12	HYP	CB-CG-CD	3.45	107.41	103.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	6	HYP	CB-CG-CD	3.52	107.49	103.14
2	J	27	HYP	CB-CG-CD	3.59	107.58	103.14
2	G	12	HYP	CB-CG-CD	3.59	107.58	103.14

All (47) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	H	3	HYP	CG
2	J	24	HYP	CG
2	H	6	HYP	CG
2	E	12	HYP	CG
2	G	27	HYP	CG
2	F	24	HYP	CG
2	I	21	HYP	CG
2	F	21	HYP	CG
2	I	30	HYP	CG
2	F	6	HYP	CG
2	E	21	HYP	CG
2	E	9	HYP	CG
2	H	27	HYP	CG
2	J	12	HYP	CG
2	F	12	HYP	CG
2	I	9	HYP	CG
2	E	30	HYP	CG
2	F	3	HYP	CG
2	G	6	HYP	CG
2	I	12	HYP	CG
2	E	3	HYP	CG
2	F	30	HYP	CG
2	H	30	HYP	CG
2	E	27	HYP	CG
2	G	12	HYP	CG
2	H	24	HYP	CG
2	E	6	HYP	CG
2	I	27	HYP	CG
2	G	9	HYP	CG
2	H	12	HYP	CG
2	J	27	HYP	CG
2	H	9	HYP	CG
2	G	24	HYP	CG
2	E	24	HYP	CG
2	H	21	HYP	CG

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Mol	Chain	Res	Type	Atom
2	J	3	HYP	CG
2	F	27	HYP	CG
2	J	21	HYP	CG
2	I	6	HYP	CG
2	J	6	HYP	CG
2	G	30	HYP	CG
2	I	3	HYP	CG
2	G	21	HYP	CG
2	I	24	HYP	CG
2	F	9	HYP	CG
2	J	9	HYP	CG
2	G	3	HYP	CG

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	30	HYP	1	0
2	G	27	HYP	1	0
2	G	3	HYP	1	0
2	G	30	HYP	3	0
2	H	3	HYP	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates 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	303/303 (100%)	0.40	9 (2%) 54 47	22, 51, 98, 122	0
1	B	299/303 (98%)	0.72	37 (12%) 5 4	30, 63, 113, 151	0
1	C	299/303 (98%)	0.53	15 (5%) 32 26	21, 60, 104, 136	0
1	D	299/303 (98%)	0.57	17 (5%) 27 22	26, 56, 118, 156	0
2	E	22/30 (73%)	0.90	3 (13%) 4 3	25, 52, 107, 161	0
2	F	22/30 (73%)	0.60	2 (9%) 11 9	31, 53, 116, 154	0
2	G	22/30 (73%)	0.60	2 (9%) 11 9	36, 54, 103, 162	0
2	H	22/30 (73%)	1.19	3 (13%) 4 3	32, 66, 149, 164	0
2	I	22/30 (73%)	0.77	2 (9%) 11 9	36, 55, 124, 158	0
2	J	21/30 (70%)	0.64	2 (9%) 10 9	25, 55, 156, 164	0
All	All	1331/1392 (95%)	0.58	92 (6%) 20 16	21, 58, 113, 164	0

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	1	GLY	8.8
2	H	2	PRO	8.3
2	I	2	PRO	7.6
2	E	2	PRO	7.4
2	H	1	GLY	6.5
1	D	328	THR	6.3
2	I	1	GLY	5.9
2	F	1	GLY	5.9
1	D	167	ALA	5.8
2	F	2	PRO	5.7
1	B	327	GLY	5.5
2	J	1	GLY	4.9
1	B	304	LYS	4.9

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Mol	Chain	Res	Type	RSRZ
2	G	1	GLY	4.9
2	J	2	PRO	4.7
1	D	101	VAL	4.7
1	B	329	VAL	4.7
1	B	68	GLN	4.6
1	B	120	GLU	4.0
1	D	327	GLY	4.0
1	B	101	VAL	3.9
1	D	122	LEU	3.7
1	B	71	ASP	3.6
1	D	104	ALA	3.5
1	A	101	VAL	3.2
1	A	42	THR	3.2
1	B	69	ASN	3.1
1	C	286	GLU	3.1
1	C	207	ILE	3.0
1	C	201	VAL	3.0
1	D	168	GLY	3.0
1	B	121	LYS	3.0
2	H	4	GLY	3.0
1	B	64	ASN	2.9
1	C	308	ASN	2.8
1	A	325	ILE	2.8
1	D	329	VAL	2.8
1	B	97	LYS	2.8
1	D	141	THR	2.8
1	C	303	GLY	2.8
1	B	175	TYR	2.7
1	B	125	VAL	2.7
1	B	122	LEU	2.7
1	B	305	GLU	2.7
1	C	270	GLY	2.7
1	D	97	LYS	2.7
1	B	119	VAL	2.7
1	B	325	ILE	2.7
1	C	100	GLN	2.7
1	B	72	MET	2.6
1	C	217	LEU	2.6
1	D	139	THR	2.6
1	C	192	LEU	2.6
1	B	67	ILE	2.6
1	B	117	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	31	ARG	2.5
1	B	61	ASP	2.5
2	G	18	THR	2.4
1	D	105	VAL	2.4
1	C	325	ILE	2.3
1	B	94	LEU	2.3
1	B	73	ILE	2.3
1	B	205	ILE	2.3
1	B	300	GLN	2.3
1	B	156	LYS	2.2
1	B	217	LEU	2.2
1	A	329	VAL	2.2
1	B	102	GLY	2.2
1	B	328	THR	2.2
1	B	70	GLY	2.2
1	D	115	PHE	2.2
2	E	4	GLY	2.2
1	D	67	ILE	2.2
1	B	84	LYS	2.2
1	A	228	GLY	2.2
1	C	237	GLN	2.2
1	B	65	GLY	2.2
1	D	295	SER	2.2
1	A	97	LYS	2.2
1	C	66	LYS	2.2
1	B	297	ALA	2.1
1	B	126	SER	2.1
1	A	327	GLY	2.1
1	D	118	LYS	2.1
1	B	118	LYS	2.1
1	A	94	LEU	2.0
1	C	219	LEU	2.0
1	C	260	THR	2.0
1	B	115	PHE	2.0
1	D	164	LYS	2.0
1	A	102	GLY	2.0
1	C	103	GLN	2.0

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

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. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	HYP	H	3	8/9	0.64	0.67	-	73,79,79,101	1
2	HYP	J	24	8/9	0.93	0.22	-	53,55,56,57	0
2	HYP	H	6	8/9	0.85	0.17	-	88,99,110,110	0
2	HYP	E	12	8/9	0.94	0.28	-	31,32,36,48	0
2	HYP	G	27	8/9	0.86	0.32	-	55,59,62,64	0
2	HYP	F	24	8/9	0.92	0.29	-	47,48,56,57	0
2	HYP	I	21	8/9	0.91	0.23	-	25,36,39,40	0
2	HYP	F	21	8/9	0.91	0.27	-	45,48,53,53	0
2	HYP	I	30	8/9	0.78	0.32	-	79,90,92,92	0
2	HYP	F	6	8/9	0.92	0.28	-	78,85,135,135	0
2	HYP	E	21	8/9	0.92	0.25	-	41,42,45,61	0
2	HYP	E	9	8/9	0.94	0.21	-	40,43,46,50	0
2	HYP	H	27	8/9	0.90	0.30	-	50,54,61,63	0
2	HYP	J	12	8/9	0.90	0.26	-	44,47,50,61	0
2	HYP	F	12	8/9	0.92	0.26	-	42,48,53,54	0
2	HYP	I	9	8/9	0.91	0.27	-	46,51,52,53	0
2	HYP	E	30	8/9	0.83	0.21	-	59,61,64,65	0
2	HYP	F	3	8/9	0.83	0.51	-	79,79,90,90	1
2	HYP	G	6	8/9	0.88	0.29	-	39,49,50,51	0
2	HYP	I	12	8/9	0.89	0.30	-	41,52,55,58	0
2	HYP	E	3	8/9	0.69	0.74	-	79,79,90,97	0
2	HYP	F	30	8/9	0.74	0.45	-	75,99,104,105	0
2	HYP	H	30	8/9	0.88	0.26	-	80,101,156,156	0
2	HYP	E	27	8/9	0.90	0.26	-	57,68,91,91	0
2	HYP	G	12	8/9	0.95	0.18	-	35,40,43,43	0
2	HYP	H	24	8/9	0.91	0.25	-	32,38,43,58	0
2	HYP	E	6	8/9	0.88	0.24	-	83,94,98,99	0
2	HYP	I	27	8/9	0.91	0.22	-	61,69,73,76	0
2	HYP	G	9	8/9	0.95	0.31	-	37,50,81,81	0
2	HYP	H	12	8/9	0.93	0.26	-	24,31,33,35	0
2	HYP	J	27	8/9	0.81	0.30	-	50,73,83,84	0
2	HYP	H	9	8/9	0.94	0.26	-	36,42,45,46	0
2	HYP	G	24	8/9	0.90	0.25	-	41,47,52,52	0
2	HYP	E	24	8/9	0.89	0.25	-	42,46,47,49	0
2	HYP	H	21	8/9	0.88	0.36	-	41,57,60,63	0
2	HYP	J	3	8/9	0.78	0.47	-	94,110,119,122	0
2	HYP	F	27	8/9	0.90	0.30	-	55,58,62,69	0
2	HYP	J	21	8/9	0.91	0.31	-	35,42,46,54	0
2	HYP	I	6	8/9	0.92	0.22	-	68,76,82,82	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	HYP	J	6	8/9	0.93	0.30	-	37,45,50,50	0
2	HYP	G	30	8/9	0.71	0.55	-	49,57,64,65	0
2	HYP	I	3	8/9	0.58	0.83	-	73,79,79,91	0
2	HYP	G	21	8/9	0.95	0.28	-	41,46,54,60	0
2	HYP	I	24	8/9	0.92	0.29	-	53,56,58,59	0
2	HYP	F	9	8/9	0.92	0.25	-	44,45,50,52	0
2	HYP	J	9	8/9	0.94	0.21	-	38,45,47,48	0
2	HYP	G	3	8/9	0.89	0.19	-	92,99,106,109	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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.