



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 03:05 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 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/ValidationPDFNotes.html>

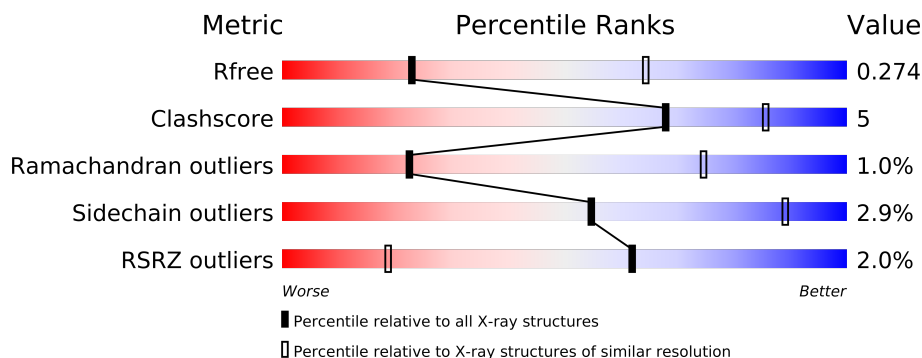
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	1341 (3.40-3.20)
Clashscore	79885	1696 (3.40-3.20)
Ramachandran outliers	78287	1664 (3.40-3.20)
Sidechain outliers	78261	1662 (3.40-3.20)
RSRZ outliers	66119	1342 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	303	
1	B	303	
1	C	303	
1	D	303	
2	E	30	
2	F	30	
2	G	30	
2	H	30	
2	I	30	
2	J	30	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10394 atoms, of which 0 are hydrogen and 0 are deuterium.

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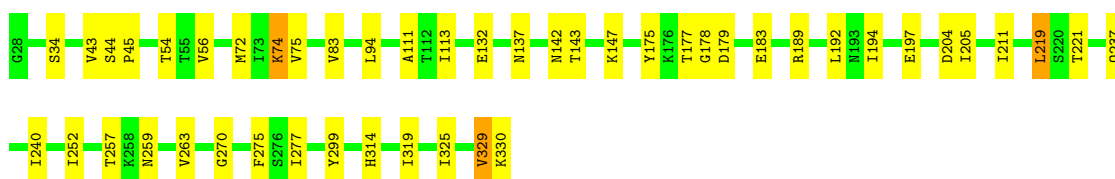
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	I	30	Total	C	N	O	0	0	0
			196	121	36	39			
2	J	28	Total	C	N	O	0	0	0
			181	111	34	36			

3 Residue-property plots

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.

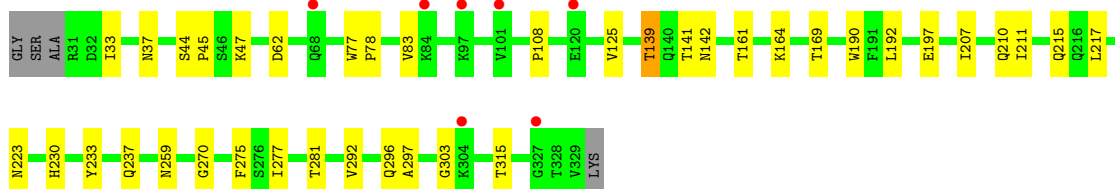
• Molecule 1: Collagen adhesin

Chain A:



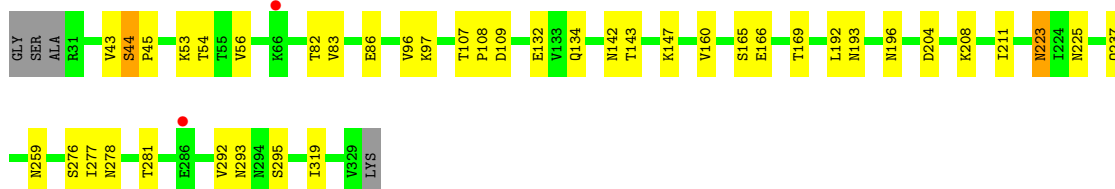
• Molecule 1: Collagen adhesin

Chain B:



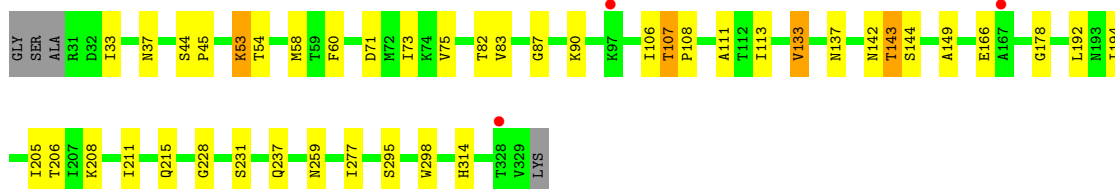
• Molecule 1: Collagen adhesin

Chain C:



• Molecule 1: Collagen adhesin

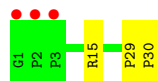
Chain D:



● Molecule 2: Collagen

Chain E: 

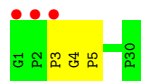
● Molecule 2: Collagen

Chain F: 

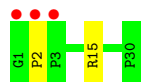
● Molecule 2: Collagen

Chain G: 

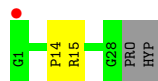
● Molecule 2: Collagen

Chain H: 

● Molecule 2: Collagen

Chain I: 

● Molecule 2: Collagen

Chain J: 

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 0.277 , 0.274	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)	0.28 , 9.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	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.

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 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

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

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 5.

All (99) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(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	Distance(Å)	Clash(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
1:B:139:THR:O	1:B:164:LYS:HG2	2.14	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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
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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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

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	301/303 (99%)	279 (93%)	21 (7%)	1 (0%)	50	92
1	B	297/303 (98%)	277 (93%)	16 (5%)	4 (1%)	18	69
1	C	297/303 (98%)	276 (93%)	16 (5%)	5 (2%)	14	62
1	D	297/303 (98%)	275 (93%)	20 (7%)	2 (1%)	30	81
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%)	4	30
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%)	22	74

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	142	ASN
1	D	142	ASN

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Mol	Chain	Res	Type
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%)	35	78
1	B	262/265 (99%)	259 (99%)	3 (1%)	84	96
1	C	262/265 (99%)	253 (97%)	9 (3%)	49	87
1	D	262/265 (99%)	257 (98%)	5 (2%)	69	93
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%)	16	56
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%)	14	49
All	All	1121/1132 (99%)	1089 (97%)	32 (3%)	55	89

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

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Mol	Chain	Res	Type
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
1	A	296	GLN
1	A	314	HIS

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Mol	Chain	Res	Type
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 chains 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	8,8,9	6.43	2 (25%)	8,10,12	1.35	2 (25%)
2	HYP	E	21	2	8,8,9	6.51	2 (25%)	8,10,12	1.62	3 (37%)
2	HYP	E	24	2	8,8,9	6.47	2 (25%)	8,10,12	1.18	1 (12%)
2	HYP	E	27	2	8,8,9	6.51	2 (25%)	8,10,12	1.12	0
2	HYP	E	3	2	8,8,9	6.42	2 (25%)	8,10,12	1.61	3 (37%)
2	HYP	E	30	2	8,8,9	6.33	2 (25%)	8,10,12	1.47	2 (25%)
2	HYP	E	6	2	8,8,9	6.50	2 (25%)	8,10,12	1.67	2 (25%)
2	HYP	E	9	2	8,8,9	6.50	2 (25%)	8,10,12	1.22	0
2	HYP	F	12	2	8,8,9	6.51	2 (25%)	8,10,12	1.43	2 (25%)
2	HYP	F	21	2	8,8,9	6.43	2 (25%)	8,10,12	1.54	3 (37%)
2	HYP	F	24	2	8,8,9	6.45	2 (25%)	8,10,12	1.59	3 (37%)
2	HYP	F	27	2	8,8,9	6.45	2 (25%)	8,10,12	1.34	1 (12%)
2	HYP	F	3	2	8,8,9	6.51	2 (25%)	8,10,12	1.52	3 (37%)
2	HYP	F	30	2	8,8,9	6.39	2 (25%)	8,10,12	1.56	3 (37%)
2	HYP	F	6	2	8,8,9	6.48	2 (25%)	8,10,12	1.56	3 (37%)
2	HYP	F	9	2	8,8,9	6.51	2 (25%)	8,10,12	1.49	3 (37%)
2	HYP	G	12	2	8,8,9	6.51	2 (25%)	8,10,12	1.60	3 (37%)
2	HYP	G	21	2	8,8,9	6.51	2 (25%)	8,10,12	1.46	3 (37%)
2	HYP	G	24	2	8,8,9	6.47	2 (25%)	8,10,12	1.59	3 (37%)
2	HYP	G	27	2	8,8,9	6.44	2 (25%)	8,10,12	1.35	1 (12%)
2	HYP	G	3	2	8,8,9	6.44	2 (25%)	8,10,12	1.48	2 (25%)
2	HYP	G	30	2	8,8,9	6.36	2 (25%)	8,10,12	1.26	1 (12%)
2	HYP	G	6	2	8,8,9	6.51	2 (25%)	8,10,12	1.48	3 (37%)
2	HYP	G	9	2	8,8,9	6.49	2 (25%)	8,10,12	1.54	3 (37%)
2	HYP	H	12	2	8,8,9	6.47	2 (25%)	8,10,12	1.19	0
2	HYP	H	21	2	8,8,9	6.50	2 (25%)	8,10,12	1.46	3 (37%)
2	HYP	H	24	2	8,8,9	6.43	2 (25%)	8,10,12	1.52	2 (25%)
2	HYP	H	27	2	8,8,9	6.54	2 (25%)	8,10,12	1.37	3 (37%)
2	HYP	H	3	2	8,8,9	6.47	2 (25%)	8,10,12	1.48	2 (25%)
2	HYP	H	30	2	8,8,9	6.36	2 (25%)	8,10,12	1.59	3 (37%)
2	HYP	H	6	2	8,8,9	6.49	2 (25%)	8,10,12	1.45	2 (25%)
2	HYP	H	9	2	8,8,9	6.44	2 (25%)	8,10,12	1.26	1 (12%)
2	HYP	I	12	2	8,8,9	6.49	2 (25%)	8,10,12	1.52	3 (37%)
2	HYP	I	21	2	8,8,9	6.47	2 (25%)	8,10,12	1.40	2 (25%)
2	HYP	I	24	2	8,8,9	6.47	2 (25%)	8,10,12	1.57	2 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HYP	I	27	2	8,8,9	6.45	2 (25%)	8,10,12	1.26	1 (12%)
2	HYP	I	3	2	8,8,9	6.48	2 (25%)	8,10,12	1.76	3 (37%)
2	HYP	I	30	2	8,8,9	6.37	2 (25%)	8,10,12	1.62	3 (37%)
2	HYP	I	6	2	8,8,9	6.46	2 (25%)	8,10,12	1.46	2 (25%)
2	HYP	I	9	2	8,8,9	6.49	2 (25%)	8,10,12	1.60	3 (37%)
2	HYP	J	12	2	8,8,9	6.51	2 (25%)	8,10,12	1.39	2 (25%)
2	HYP	J	21	2	8,8,9	6.50	2 (25%)	8,10,12	1.51	3 (37%)
2	HYP	J	24	2	8,8,9	6.52	2 (25%)	8,10,12	1.56	3 (37%)
2	HYP	J	27	2	8,8,9	6.52	2 (25%)	8,10,12	1.62	3 (37%)
2	HYP	J	3	2	8,8,9	6.50	2 (25%)	8,10,12	1.50	3 (37%)
2	HYP	J	6	2	8,8,9	6.50	2 (25%)	8,10,12	1.58	3 (37%)
2	HYP	J	9	2	8,8,9	6.48	2 (25%)	8,10,12	1.57	3 (37%)

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

All (94) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	27	HYP	O-C	18.22	1.23	1.11
2	G	12	HYP	O-C	18.17	1.23	1.11
2	J	27	HYP	O-C	18.17	1.23	1.11
2	J	24	HYP	O-C	18.16	1.23	1.11
2	F	3	HYP	O-C	18.15	1.23	1.11
2	F	12	HYP	O-C	18.15	1.23	1.11
2	G	21	HYP	O-C	18.15	1.23	1.11
2	F	9	HYP	O-C	18.15	1.23	1.11
2	G	6	HYP	O-C	18.15	1.23	1.11
2	E	27	HYP	O-C	18.14	1.23	1.11
2	J	12	HYP	O-C	18.14	1.23	1.11
2	E	21	HYP	O-C	18.13	1.23	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	6	HYP	O-C	18.13	1.23	1.11
2	J	21	HYP	O-C	18.12	1.23	1.11
2	J	3	HYP	O-C	18.11	1.23	1.11
2	H	21	HYP	O-C	18.11	1.23	1.11
2	E	6	HYP	O-C	18.09	1.23	1.11
2	E	9	HYP	O-C	18.09	1.23	1.11
2	G	9	HYP	O-C	18.09	1.23	1.11
2	J	9	HYP	O-C	18.08	1.23	1.11
2	H	6	HYP	O-C	18.08	1.23	1.11
2	I	9	HYP	O-C	18.07	1.23	1.11
2	I	12	HYP	O-C	18.07	1.23	1.11
2	H	3	HYP	O-C	18.04	1.23	1.11
2	F	6	HYP	O-C	18.03	1.23	1.11
2	G	24	HYP	O-C	18.03	1.23	1.11
2	H	12	HYP	O-C	18.03	1.23	1.11
2	I	3	HYP	O-C	18.02	1.23	1.11
2	I	21	HYP	O-C	18.02	1.23	1.11
2	I	24	HYP	O-C	18.01	1.23	1.11
2	E	24	HYP	O-C	18.00	1.23	1.11
2	I	6	HYP	O-C	17.99	1.23	1.11
2	I	27	HYP	O-C	17.97	1.23	1.11
2	F	24	HYP	O-C	17.96	1.23	1.11
2	G	3	HYP	O-C	17.95	1.23	1.11
2	H	9	HYP	O-C	17.95	1.23	1.11
2	F	27	HYP	O-C	17.93	1.23	1.11
2	G	27	HYP	O-C	17.92	1.23	1.11
2	F	21	HYP	O-C	17.91	1.23	1.11
2	E	12	HYP	O-C	17.90	1.23	1.11
2	E	3	HYP	O-C	17.86	1.23	1.11
2	H	24	HYP	O-C	17.84	1.23	1.11
2	F	30	HYP	O-C	17.79	1.23	1.11
2	I	30	HYP	O-C	17.75	1.23	1.11
2	H	30	HYP	O-C	17.71	1.23	1.11
2	G	30	HYP	O-C	17.70	1.23	1.11
2	E	30	HYP	O-C	17.61	1.23	1.11
2	H	24	HYP	CA-C	3.24	1.54	1.48
2	E	6	HYP	CA-C	3.07	1.54	1.48
2	E	3	HYP	CA-C	3.07	1.54	1.48
2	I	3	HYP	CA-C	3.07	1.54	1.48
2	F	27	HYP	CA-C	3.02	1.54	1.48
2	E	12	HYP	CA-C	3.00	1.53	1.48
2	H	6	HYP	CA-C	2.97	1.53	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	24	HYP	CA-C	2.97	1.53	1.48
2	G	27	HYP	CA-C	2.97	1.53	1.48
2	E	9	HYP	CA-C	2.96	1.53	1.48
2	E	27	HYP	CA-C	2.95	1.53	1.48
2	H	12	HYP	CA-C	2.94	1.53	1.48
2	H	27	HYP	CA-C	2.94	1.53	1.48
2	F	6	HYP	CA-C	2.93	1.53	1.48
2	E	24	HYP	CA-C	2.93	1.53	1.48
2	F	24	HYP	CA-C	2.92	1.53	1.48
2	E	30	HYP	CA-C	2.91	1.53	1.48
2	J	27	HYP	CA-C	2.91	1.53	1.48
2	H	21	HYP	CA-C	2.91	1.53	1.48
2	I	6	HYP	CA-C	2.90	1.53	1.48
2	G	30	HYP	CA-C	2.90	1.53	1.48
2	I	12	HYP	CA-C	2.90	1.53	1.48
2	F	9	HYP	CA-C	2.90	1.53	1.48
2	I	9	HYP	CA-C	2.89	1.53	1.48
2	E	21	HYP	CA-C	2.89	1.53	1.48
2	G	24	HYP	CA-C	2.89	1.53	1.48
2	F	21	HYP	CA-C	2.89	1.53	1.48
2	F	30	HYP	CA-C	2.89	1.53	1.48
2	H	9	HYP	CA-C	2.88	1.53	1.48
2	F	12	HYP	CA-C	2.87	1.53	1.48
2	I	30	HYP	CA-C	2.87	1.53	1.48
2	J	24	HYP	CA-C	2.85	1.53	1.48
2	H	30	HYP	CA-C	2.85	1.53	1.48
2	J	3	HYP	CA-C	2.85	1.53	1.48
2	J	12	HYP	CA-C	2.85	1.53	1.48
2	G	9	HYP	CA-C	2.84	1.53	1.48
2	F	3	HYP	CA-C	2.83	1.53	1.48
2	I	27	HYP	CA-C	2.83	1.53	1.48
2	H	3	HYP	CA-C	2.82	1.53	1.48
2	J	21	HYP	CA-C	2.82	1.53	1.48
2	G	3	HYP	CA-C	2.82	1.53	1.48
2	I	21	HYP	CA-C	2.82	1.53	1.48
2	G	21	HYP	CA-C	2.82	1.53	1.48
2	G	6	HYP	CA-C	2.81	1.53	1.48
2	J	9	HYP	CA-C	2.81	1.53	1.48
2	J	6	HYP	CA-C	2.77	1.53	1.48
2	G	12	HYP	CA-C	2.74	1.53	1.48

All (108) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	3	HYP	CG-CB-CA	2.97	108.55	104.22
2	G	12	HYP	CB-CG-CD	2.90	107.58	103.46
2	J	27	HYP	CB-CG-CD	2.90	107.58	103.46
2	H	24	HYP	C-CA-N	2.88	116.18	110.86
2	J	6	HYP	CB-CG-CD	2.83	107.49	103.46
2	J	12	HYP	CB-CG-CD	2.77	107.41	103.46
2	E	21	HYP	CB-CG-CD	2.76	107.39	103.46
2	H	30	HYP	CG-CB-CA	2.76	108.25	104.22
2	E	6	HYP	CG-CB-CA	2.75	108.23	104.22
2	I	30	HYP	CB-CG-CD	2.74	107.36	103.46
2	F	3	HYP	CB-CG-CD	2.73	107.35	103.46
2	I	30	HYP	CG-CB-CA	2.70	108.17	104.22
2	J	9	HYP	CB-CG-CD	2.70	107.30	103.46
2	I	12	HYP	CB-CG-CD	2.69	107.29	103.46
2	I	24	HYP	CB-CG-CD	2.68	107.28	103.46
2	F	30	HYP	CB-CG-CD	2.67	107.26	103.46
2	I	9	HYP	CB-CG-CD	2.67	107.26	103.46
2	G	24	HYP	CB-CG-CD	2.67	107.25	103.46
2	F	30	HYP	CG-CB-CA	2.66	108.11	104.22
2	E	6	HYP	CB-CG-CD	2.65	107.23	103.46
2	J	24	HYP	CG-CB-CA	2.65	108.08	104.22
2	F	24	HYP	CB-CG-CD	2.64	107.22	103.46
2	F	6	HYP	CG-CB-CA	2.63	108.06	104.22
2	F	24	HYP	CG-CB-CA	2.62	108.05	104.22
2	F	6	HYP	CB-CG-CD	2.62	107.18	103.46
2	J	24	HYP	CB-CG-CD	2.61	107.18	103.46
2	J	9	HYP	CG-CB-CA	2.61	108.04	104.22
2	G	24	HYP	CG-CB-CA	2.61	108.03	104.22
2	J	21	HYP	CB-CG-CD	2.60	107.16	103.46
2	G	9	HYP	CG-CB-CA	2.60	108.02	104.22
2	H	21	HYP	CB-CG-CD	2.59	107.14	103.46
2	E	30	HYP	CG-CB-CA	2.58	107.99	104.22
2	E	21	HYP	CG-CB-CA	2.58	107.99	104.22
2	F	9	HYP	CB-CG-CD	2.58	107.13	103.46
2	F	21	HYP	CB-CG-CD	2.58	107.12	103.46
2	G	9	HYP	CB-CG-CD	2.57	107.11	103.46
2	I	9	HYP	CG-CB-CA	2.56	107.95	104.22
2	G	21	HYP	CB-CG-CD	2.56	107.10	103.46
2	H	3	HYP	CG-CB-CA	2.55	107.94	104.22
2	I	3	HYP	C-CA-N	2.54	115.57	110.86
2	G	3	HYP	CG-CB-CA	2.54	107.93	104.22
2	G	6	HYP	CB-CG-CD	2.54	107.07	103.46
2	J	3	HYP	CB-CG-CD	2.53	107.07	103.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	3	HYP	CB-CG-CD	2.53	107.06	103.46
2	F	21	HYP	CG-CB-CA	2.53	107.92	104.22
2	I	6	HYP	CG-CB-CA	2.53	107.91	104.22
2	E	3	HYP	CB-CG-CD	2.52	107.05	103.46
2	G	3	HYP	CB-CG-CD	2.52	107.05	103.46
2	F	12	HYP	CG-CB-CA	2.52	107.90	104.22
2	I	24	HYP	CG-CB-CA	2.51	107.89	104.22
2	I	6	HYP	CB-CG-CD	2.47	106.98	103.46
2	F	3	HYP	CG-CB-CA	2.45	107.80	104.22
2	E	3	HYP	CG-CB-CA	2.44	107.78	104.22
2	H	30	HYP	CB-CG-CD	2.43	106.92	103.46
2	J	3	HYP	CG-CB-CA	2.42	107.76	104.22
2	H	24	HYP	CB-CG-CD	2.42	106.91	103.46
2	G	6	HYP	CG-CB-CA	2.42	107.75	104.22
2	F	12	HYP	CB-CG-CD	2.41	106.89	103.46
2	H	6	HYP	CG-CB-CA	2.41	107.74	104.22
2	F	9	HYP	CG-CB-CA	2.41	107.73	104.22
2	E	12	HYP	CB-CG-CD	2.37	106.83	103.46
2	H	6	HYP	CB-CG-CD	2.36	106.81	103.46
2	I	12	HYP	CG-CB-CA	2.35	107.66	104.22
2	I	21	HYP	CB-CG-CD	2.34	106.80	103.46
2	G	21	HYP	CD-N-CA	2.33	112.54	108.15
2	H	27	HYP	CB-CG-CD	2.32	106.75	103.46
2	J	6	HYP	CG-CB-CA	2.32	107.60	104.22
2	J	6	HYP	CD-N-CA	2.31	112.51	108.15
2	J	27	HYP	C-CA-N	2.30	115.11	110.86
2	I	3	HYP	CB-CG-CD	2.29	106.72	103.46
2	J	21	HYP	CD-N-CA	2.26	112.42	108.15
2	I	21	HYP	CG-CB-CA	2.26	107.52	104.22
2	G	12	HYP	CG-CB-CA	2.25	107.50	104.22
2	J	24	HYP	CD-N-CA	2.22	112.35	108.15
2	G	30	HYP	CG-CB-CA	2.22	107.46	104.22
2	J	9	HYP	CD-N-CA	2.22	112.34	108.15
2	J	21	HYP	CG-CB-CA	2.21	107.44	104.22
2	G	24	HYP	CD-N-CA	2.20	112.31	108.15
2	E	30	HYP	CB-CG-CD	2.20	106.59	103.46
2	I	30	HYP	CD-N-CA	2.19	112.29	108.15
2	G	12	HYP	CD-N-CA	2.16	112.23	108.15
2	G	9	HYP	CD-N-CA	2.16	112.22	108.15
2	H	27	HYP	CD-N-CA	2.14	112.19	108.15
2	I	9	HYP	CD-N-CA	2.13	112.18	108.15
2	F	21	HYP	CD-N-CA	2.13	112.17	108.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	9	HYP	CD-N-CA	2.13	112.17	108.15
2	H	30	HYP	CD-N-CA	2.13	112.16	108.15
2	I	12	HYP	CD-N-CA	2.12	112.16	108.15
2	G	27	HYP	CG-CB-CA	2.10	107.29	104.22
2	G	6	HYP	CD-N-CA	2.10	112.11	108.15
2	J	12	HYP	CG-CB-CA	2.10	107.28	104.22
2	F	3	HYP	CD-N-CA	2.09	112.10	108.15
2	F	30	HYP	CD-N-CA	2.09	112.09	108.15
2	J	27	HYP	CG-CB-CA	2.09	107.27	104.22
2	H	21	HYP	CD-N-CA	2.09	112.09	108.15
2	J	3	HYP	CD-N-CA	2.09	112.09	108.15
2	E	21	HYP	CD-N-CA	2.08	112.07	108.15
2	F	6	HYP	CD-N-CA	2.07	112.07	108.15
2	E	3	HYP	CD-N-CA	2.07	112.05	108.15
2	H	21	HYP	CG-CB-CA	2.06	107.23	104.22
2	E	12	HYP	CD-N-CA	2.05	112.02	108.15
2	G	21	HYP	CG-CB-CA	2.05	107.21	104.22
2	F	27	HYP	CB-CG-CD	2.04	106.37	103.46
2	H	9	HYP	CG-CB-CA	2.03	107.19	104.22
2	I	27	HYP	CB-CG-CD	2.02	106.33	103.46
2	E	24	HYP	CB-CG-CD	2.01	106.33	103.46
2	F	24	HYP	CD-N-CA	2.01	111.94	108.15
2	H	27	HYP	CG-CB-CA	2.00	107.14	104.22

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

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Mol	Chain	Res	Type	Atom
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
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.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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	303/303 (100%)	0.16	0 100 100	22, 51, 98, 122	0
1	B	299/303 (98%)	0.39	7 (2%) 57 15	30, 63, 113, 151	0
1	C	299/303 (98%)	0.29	2 (0%) 84 42	21, 60, 104, 136	0
1	D	299/303 (98%)	0.20	3 (1%) 79 33	26, 56, 118, 156	0
2	E	30/30 (100%)	0.47	3 (10%) 8 2	25, 52, 107, 161	0
2	F	30/30 (100%)	0.43	3 (10%) 8 2	31, 53, 116, 154	0
2	G	30/30 (100%)	0.25	2 (6%) 17 4	36, 54, 103, 162	0
2	H	30/30 (100%)	0.59	3 (10%) 8 2	31, 64, 149, 164	0
2	I	30/30 (100%)	0.45	3 (10%) 8 2	36, 56, 124, 158	0
2	J	28/30 (93%)	0.34	1 (3%) 41 9	25, 55, 156, 164	0
All	All	1378/1392 (98%)	0.28	27 (1%) 62 19	21, 58, 112, 164	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	3	HYP	5.1
2	H	2	PRO	4.3
2	I	3	HYP	4.1
2	E	1	GLY	4.1
2	I	2	PRO	3.9
2	E	2	PRO	3.8
2	F	2	PRO	3.8
2	H	3	HYP	3.7
2	H	1	GLY	3.6
2	F	1	GLY	3.6
1	D	167	ALA	3.2
1	D	328	THR	3.0
1	B	120	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	304	LYS	2.5
1	B	97	LYS	2.4
2	G	1	GLY	2.4
2	I	1	GLY	2.4
1	B	68	GLN	2.4
2	J	1	GLY	2.3
2	F	3	HYP	2.2
1	C	66	LYS	2.2
1	B	101	VAL	2.1
1	C	286	GLU	2.1
1	D	97	LYS	2.1
1	B	327	GLY	2.1
2	G	30	HYP	2.0
1	B	84	LYS	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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	HYP	F	30	8/9	0.47	4.47	75,99,104,105	0
2	HYP	G	30	8/9	0.50	3.40	49,57,64,65	0
2	HYP	G	9	8/9	0.32	2.32	37,50,81,81	0
2	HYP	J	27	8/9	0.32	1.76	50,73,83,84	0
2	HYP	G	6	8/9	0.29	1.68	39,49,50,51	0
2	HYP	H	21	8/9	0.36	1.66	41,57,60,63	0
2	HYP	I	30	8/9	0.30	1.57	79,90,92,92	0
2	HYP	J	21	8/9	0.34	1.54	35,42,46,54	0
2	HYP	G	21	8/9	0.31	1.43	41,46,54,60	0
2	HYP	F	21	8/9	0.27	1.35	45,48,53,53	0
2	HYP	F	27	8/9	0.29	1.20	55,58,62,69	0
2	HYP	F	24	8/9	0.25	1.18	47,48,56,57	0
2	HYP	I	3	8/9	0.75	1.13	73,79,79,91	0
2	HYP	E	12	8/9	0.28	1.11	31,32,36,48	0
2	HYP	F	6	8/9	0.37	0.74	78,85,135,135	0
2	HYP	I	9	8/9	0.28	0.74	46,51,52,53	0
2	HYP	G	24	8/9	0.25	0.65	41,47,52,52	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	HYP	J	6	8/9	0.33	0.62	37,45,50,50	0
2	HYP	F	9	8/9	0.22	0.62	44,45,50,52	0
2	HYP	I	24	8/9	0.32	0.59	53,56,58,59	0
2	HYP	E	24	8/9	0.26	0.53	42,46,47,49	0
2	HYP	E	3	8/9	0.86	0.51	79,79,90,97	0
2	HYP	J	12	8/9	0.29	0.50	44,47,50,61	0
2	HYP	I	12	8/9	0.33	0.44	41,52,55,58	0
2	HYP	E	27	8/9	0.25	0.43	57,68,91,91	0
2	HYP	F	12	8/9	0.24	0.35	42,48,53,54	0
2	HYP	H	27	8/9	0.30	0.35	50,54,61,63	0
2	HYP	H	12	8/9	0.24	0.26	24,31,33,35	0
2	HYP	G	27	8/9	0.26	0.22	55,59,62,64	0
2	HYP	J	3	8/9	0.44	0.20	94,110,119,122	0
2	HYP	H	24	8/9	0.26	0.11	32,38,43,58	0
2	HYP	I	21	8/9	0.24	0.11	25,36,39,40	0
2	HYP	F	3	8/9	0.52	0.06	79,79,90,90	1
2	HYP	E	21	8/9	0.23	-0.01	41,42,45,61	0
2	HYP	E	9	8/9	0.19	-0.06	40,43,46,50	0
2	HYP	H	9	8/9	0.27	-0.07	36,42,45,46	0
2	HYP	H	30	8/9	0.27	-0.13	80,101,156,156	0
2	HYP	H	3	8/9	0.70	-0.14	73,79,79,101	1
2	HYP	J	9	8/9	0.22	-0.16	38,45,47,48	0
2	HYP	J	24	8/9	0.24	-0.17	53,55,56,57	0
2	HYP	E	6	8/9	0.28	-0.21	83,94,98,99	0
2	HYP	I	6	8/9	0.26	-0.22	68,76,82,82	0
2	HYP	I	27	8/9	0.23	-0.24	61,69,73,76	0
2	HYP	E	30	8/9	0.20	-0.46	59,61,64,65	0
2	HYP	G	12	8/9	0.18	-0.77	35,40,43,43	0
2	HYP	G	3	8/9	0.21	-1.09	92,99,106,109	0
2	HYP	H	6	8/9	0.18	-1.34	88,99,110,110	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.