



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:46 PM GMT

PDB ID : 1YK3
Title : Crystal structure of Rv1347c from Mycobacterium tuberculosis
Authors : Card, G.L.; Peterson, N.A.; Smith, C.A.; Rupp, B.; Schick, B.M.; Baker, E.N.;
TB Structural Genomics Consortium (TBSGC)
Deposited on : 2005-01-16
Resolution : 2.20 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

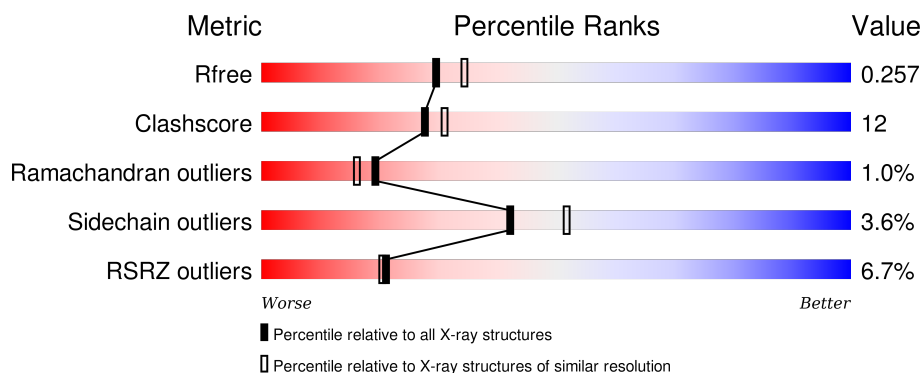
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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	210	
1	B	210	
1	C	210	
1	D	210	
1	E	210	

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Mol	Chain	Length	Quality of chain
1	F	210	
1	G	210	
1	H	210	

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	BOG	D	701	-	-	-	X
2	BOG	E	702	-	-	-	X
2	BOG	G	703	-	-	-	X

2 Entry composition

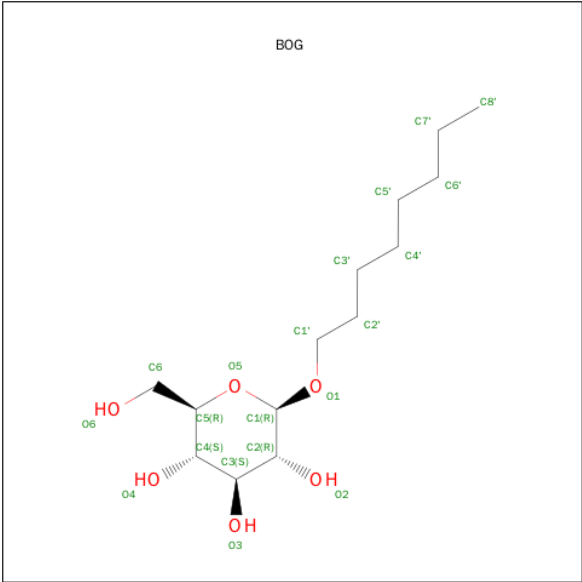
There are 3 unique types of molecules in this entry. The entry contains 13515 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Hypothetical protein Rv1347c/MT1389.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	198	Total	C	N	O	S	0	0	0
			1600	1017	292	284	7			
1	B	193	Total	C	N	O	S	0	0	0
			1563	995	285	276	7			
1	C	199	Total	C	N	O	S	0	0	0
			1607	1021	293	286	7			
1	D	200	Total	C	N	O	S	0	0	0
			1612	1024	294	287	7			
1	E	196	Total	C	N	O	S	0	0	0
			1587	1010	290	280	7			
1	F	202	Total	C	N	O	S	0	0	0
			1625	1031	297	290	7			
1	G	201	Total	C	N	O	S	0	0	0
			1618	1027	295	289	7			
1	H	197	Total	C	N	O	S	0	0	0
			1593	1013	291	282	7			

- Molecule 2 is SUGAR (B-OCTYLGLUCOSIDE) (three-letter code: BOG) (formula: C₁₄H₂₈O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			20	14	6		
2	E	1	Total	C	O	0	0
			20	14	6		
2	G	1	Total	C	O	0	0
			20	14	6		

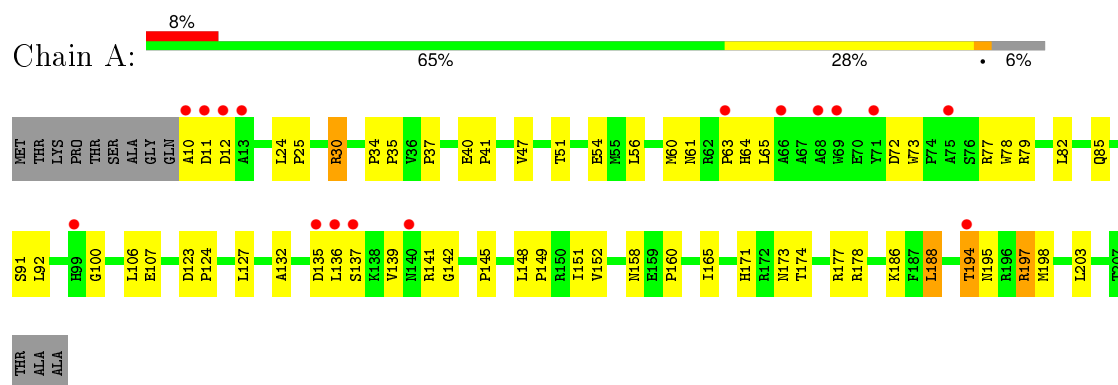
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	51	Total	O	0	0
			51	51		
3	B	54	Total	O	0	0
			54	54		
3	C	83	Total	O	0	0
			83	83		
3	D	84	Total	O	0	0
			84	84		
3	E	59	Total	O	0	0
			59	59		
3	F	136	Total	O	0	0
			136	136		
3	G	121	Total	O	0	0
			121	121		
3	H	62	Total	O	0	0
			62	62		

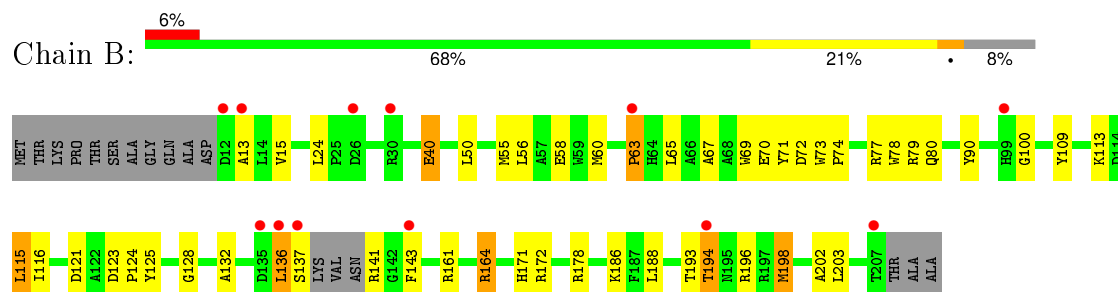
3 Residue-property plots [i](#)

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.

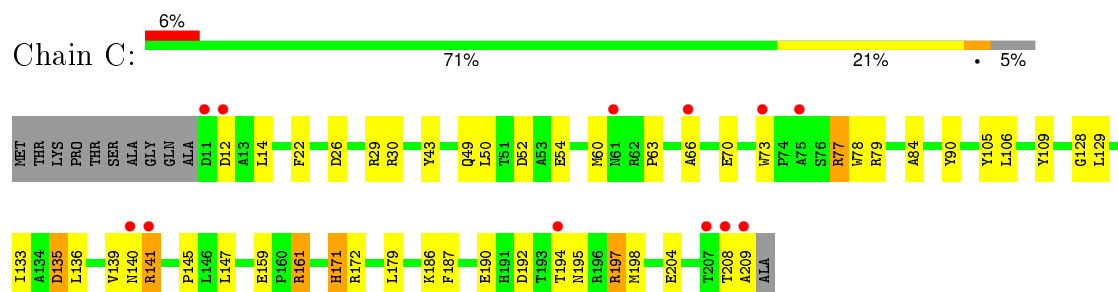
- Molecule 1: Hypothetical protein Rv1347c/MT1389



- Molecule 1: Hypothetical protein Rv1347c/MT1389

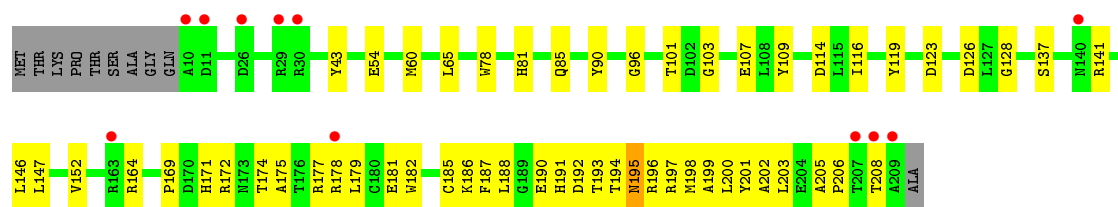


- Molecule 1: Hypothetical protein Rv1347c/MT1389

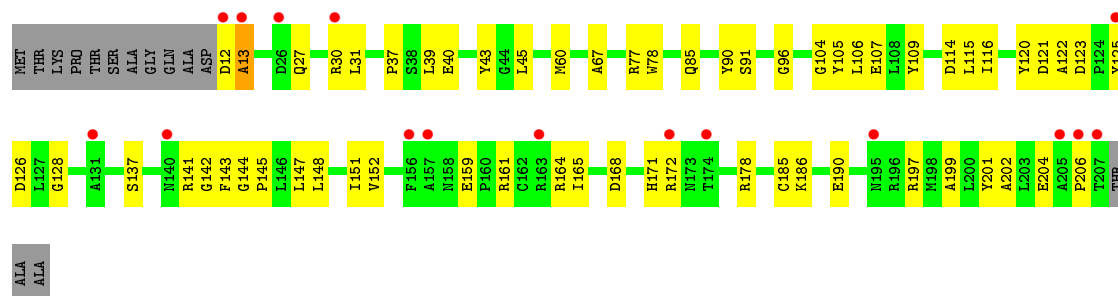


- Molecule 1: Hypothetical protein Rv1347c/MT1389

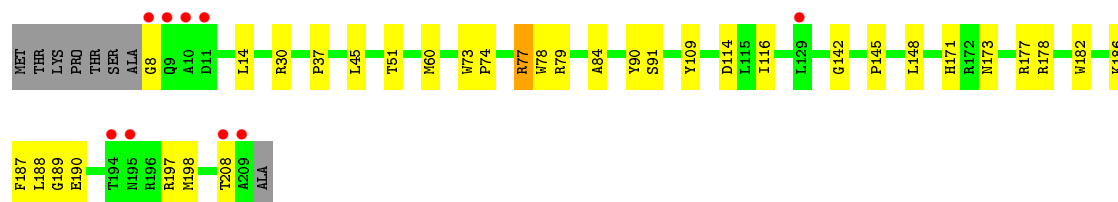
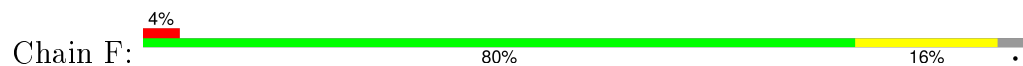




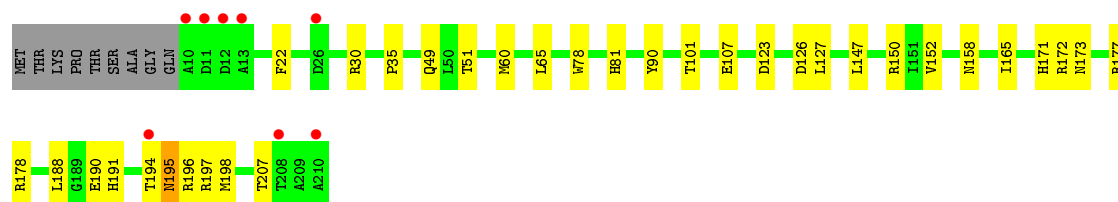
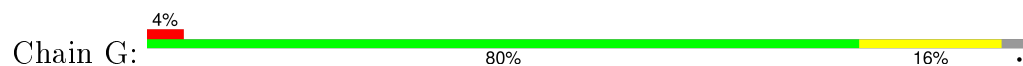
• Molecule 1: Hypothetical protein Rv1347c/MT1389



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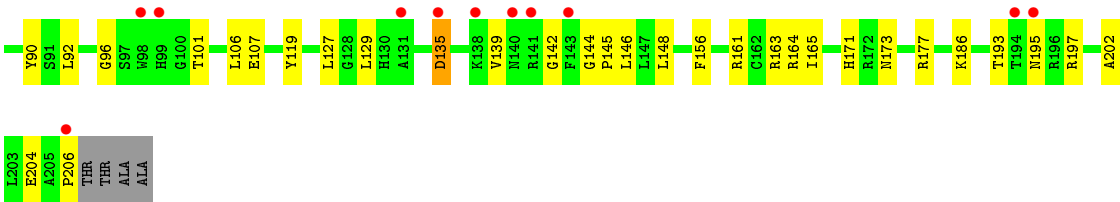


• Molecule 1: Hypothetical protein Rv1347c/MT1389



• Molecule 1: Hypothetical protein Rv1347c/MT1389





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.67Å 77.26Å 296.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.96 – 2.20 36.96 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.9 (36.96-2.20) 96.1 (36.96-2.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.81 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.227 , 0.258 0.227 , 0.257	Depositor DCC
R_{free} test set	1824 reflections (2.13%)	DCC
Wilson B-factor (Å ²)	29.4	Xtriage
Anisotropy	0.360	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 50.4	EDS
Estimated twinning fraction	0.012 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 89275 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13515	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.27% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.32	0/1649	0.57	0/2250
1	B	0.33	0/1611	0.58	0/2197
1	C	0.34	0/1656	0.62	0/2260
1	D	0.34	0/1661	0.61	0/2267
1	E	0.34	0/1636	0.61	1/2232 (0.0%)
1	F	0.38	0/1674	0.64	0/2284
1	G	0.37	0/1667	0.64	0/2274
1	H	0.32	0/1642	0.57	0/2240
All	All	0.34	0/13196	0.61	1/18004 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	199	ALA	N-CA-C	-5.03	97.42	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1600	0	1546	44	0
1	B	1563	0	1508	44	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1607	0	1553	42	0
1	D	1612	0	1558	42	0
1	E	1587	0	1537	46	0
1	F	1625	0	1569	23	0
1	G	1618	0	1563	28	0
1	H	1593	0	1539	42	0
2	D	20	0	28	3	0
2	E	20	0	28	4	0
2	G	20	0	28	2	0
3	A	51	0	0	2	0
3	B	54	0	0	6	0
3	C	83	0	0	2	0
3	D	84	0	0	1	0
3	E	59	0	0	0	0
3	F	136	0	0	2	0
3	G	121	0	0	4	0
3	H	62	0	0	3	0
All	All	13515	0	12457	298	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 12.

All (298) 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:H:60:MET:HG3	1:H:78:TRP:CZ2	2.11	0.84
1:B:60:MET:HG3	1:B:78:TRP:CZ2	2.13	0.83
1:B:136:LEU:HD23	1:B:136:LEU:H	1.46	0.80
1:F:60:MET:HG3	1:F:78:TRP:CZ2	2.18	0.79
1:H:106:LEU:HD23	1:H:129:LEU:HD11	1.66	0.77
1:D:190:GLU:HB3	1:D:197:ARG:NH1	2.00	0.77
1:A:158:ASN:HA	1:B:63:PRO:HB3	1.66	0.76
1:B:125:TYR:HE2	1:B:161:ARG:HH21	1.32	0.76
1:C:133:ILE:HB	1:C:139:VAL:HG22	1.67	0.75
1:C:60:MET:HG3	1:C:78:TRP:CZ2	2.22	0.75
1:A:10:ALA:HA	1:E:77:ARG:HH11	1.51	0.74
1:D:60:MET:HG3	1:D:78:TRP:CZ2	2.24	0.72
1:G:195:ASN:HD22	1:G:195:ASN:N	1.85	0.72
1:E:85:GLN:HE22	1:E:107:GLU:HG2	1.55	0.71
1:C:43:TYR:CE1	1:C:147:LEU:HD11	2.27	0.70
1:B:116:ILE:HD11	1:B:198:MET:HE2	1.72	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:MET:HG2	1:A:78:TRP:HZ2	1.57	0.69
1:D:172:ARG:HB2	1:D:172:ARG:NH1	2.07	0.69
1:G:60:MET:HG3	1:G:78:TRP:CZ2	2.28	0.69
1:D:190:GLU:HB3	1:D:197:ARG:HH12	1.58	0.68
1:D:96:GLY:HA3	2:D:701:BOG:H7'2	1.76	0.68
1:C:136:LEU:H	1:C:136:LEU:HD12	1.59	0.68
1:D:85:GLN:HE22	1:D:107:GLU:HG2	1.59	0.68
1:E:60:MET:HG3	1:E:78:TRP:CZ2	2.29	0.67
1:H:85:GLN:HE22	1:H:107:GLU:HG2	1.58	0.67
1:C:171:HIS:CD2	1:C:197:ARG:HB3	2.29	0.67
1:E:114:ASP:OD1	1:E:115:LEU:HD23	1.95	0.66
1:D:172:ARG:HH11	1:D:172:ARG:HB2	1.61	0.66
1:H:73:TRP:HB3	1:H:74:PRO:HD2	1.78	0.66
1:H:173:ASN:O	1:H:177:ARG:HG3	1.96	0.66
1:C:22:PHE:HB3	1:G:22:PHE:HB3	1.78	0.66
1:G:195:ASN:HD22	1:G:195:ASN:H	1.45	0.65
1:G:150:ARG:HB2	3:G:819:HOH:O	1.96	0.65
1:D:171:HIS:CD2	1:D:197:ARG:HB3	2.31	0.65
1:B:171:HIS:CE1	1:B:172:ARG:HG3	2.32	0.65
1:D:116:ILE:HD11	1:D:200:LEU:CD1	2.28	0.64
1:A:100:GLY:HA2	1:D:187:PHE:HB2	1.79	0.64
1:E:123:ASP:HB3	1:E:126:ASP:OD1	1.97	0.64
1:D:175:ALA:HA	1:D:178:ARG:HE	1.62	0.64
1:E:96:GLY:HA3	2:E:702:BOG:H7'2	1.80	0.64
1:H:56:LEU:HD12	1:H:82:LEU:HD11	1.80	0.64
1:H:171:HIS:CD2	1:H:197:ARG:HB3	2.33	0.64
1:F:8:GLY:HA3	3:F:219:HOH:O	1.97	0.63
1:A:137:SER:O	1:A:141:ARG:HD3	1.98	0.63
1:H:39:LEU:HD12	1:H:45:LEU:HG	1.81	0.63
1:E:164:ARG:HD2	1:E:202:ALA:HB1	1.81	0.63
1:D:190:GLU:CB	1:D:197:ARG:HH12	2.12	0.62
1:G:81:HIS:HD2	3:G:811:HOH:O	1.81	0.62
1:C:73:TRP:HB3	1:C:77:ARG:HB3	1.82	0.62
1:D:164:ARG:HD3	1:D:202:ALA:HB1	1.81	0.62
1:H:60:MET:CE	1:H:65:LEU:HD13	2.29	0.62
1:B:50:LEU:HD13	1:B:79:ARG:HE	1.64	0.62
1:H:142:GLY:C	1:H:145:PRO:HD2	2.20	0.62
1:B:136:LEU:H	1:B:136:LEU:CD2	2.12	0.62
1:D:174:THR:O	1:D:178:ARG:HG2	1.99	0.62
1:C:77:ARG:HG3	1:C:77:ARG:HH11	1.64	0.62
1:G:195:ASN:N	1:G:195:ASN:ND2	2.48	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:105:TYR:O	1:E:106:LEU:HD23	2.01	0.61
1:F:77:ARG:HG2	1:F:77:ARG:HH11	1.64	0.60
1:H:164:ARG:HD2	1:H:202:ALA:HB1	1.83	0.60
1:H:69:TRP:O	1:H:71:TYR:N	2.31	0.60
1:B:136:LEU:HD23	1:B:136:LEU:N	2.16	0.59
1:E:85:GLN:HG2	1:E:91:SER:HB3	1.84	0.59
1:A:37:PRO:HG3	3:A:238:HOH:O	2.03	0.59
1:E:67:ALA:HB2	1:H:161:ARG:HG2	1.84	0.59
1:C:50:LEU:HG	3:C:274:HOH:O	2.00	0.59
1:B:73:TRP:HB3	1:B:74:PRO:HD2	1.85	0.59
1:E:152:VAL:HG13	1:E:165:ILE:HD12	1.83	0.59
1:C:194:THR:O	1:C:195:ASN:HB2	2.02	0.59
1:C:135:ASP:O	1:C:139:VAL:HG23	2.03	0.58
1:E:27:GLN:HG3	1:E:30:ARG:HH21	1.68	0.58
1:D:123:ASP:HB3	1:D:126:ASP:OD1	2.02	0.58
1:B:123:ASP:OD1	1:B:124:PRO:HD2	2.03	0.58
1:H:11:ASP:OD1	1:H:79:ARG:NH1	2.36	0.58
1:H:60:MET:HE2	1:H:65:LEU:HD13	1.84	0.58
1:F:77:ARG:CG	1:F:77:ARG:HH11	2.16	0.58
1:H:135:ASP:O	1:H:139:VAL:HG23	2.03	0.57
1:B:13:ALA:HB3	3:B:225:HOH:O	2.03	0.57
1:F:171:HIS:CD2	1:F:197:ARG:HB3	2.39	0.57
1:G:196:ARG:HG2	1:G:196:ARG:HH11	1.67	0.57
1:H:58:GLU:HG2	1:H:62:ARG:NH2	2.19	0.57
1:D:101:THR:HG21	2:D:701:BOG:H2	1.87	0.57
1:C:194:THR:HG22	1:C:195:ASN:ND2	2.19	0.57
1:E:37:PRO:HG2	1:E:45:LEU:HD23	1.86	0.57
1:D:81:HIS:HE1	1:D:107:GLU:OE2	1.87	0.57
1:G:152:VAL:HG13	1:G:165:ILE:HD12	1.87	0.57
1:E:115:LEU:HD23	1:E:116:ILE:N	2.19	0.56
1:B:60:MET:CE	1:B:65:LEU:HD13	2.35	0.56
1:C:159:GLU:OE2	1:C:161:ARG:HD3	2.05	0.56
1:E:116:ILE:HD11	1:E:168:ASP:HB3	1.86	0.56
1:D:190:GLU:OE1	1:D:197:ARG:NH1	2.39	0.56
1:F:91:SER:HB3	1:F:109:TYR:HB3	1.87	0.56
1:A:135:ASP:O	1:A:139:VAL:HG23	2.06	0.55
1:A:56:LEU:HD12	1:A:82:LEU:HD11	1.88	0.55
1:G:147:LEU:HD13	2:G:703:BOG:H7'2	1.89	0.55
1:C:192:ASP:OD1	1:C:197:ARG:HD2	2.06	0.55
1:B:109:TYR:CZ	1:B:128:GLY:HA3	2.41	0.55
1:E:85:GLN:NE2	1:E:107:GLU:HG2	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:30:ARG:HH11	1:E:30:ARG:HG3	1.72	0.55
1:D:146:LEU:HD23	1:D:146:LEU:O	2.06	0.55
1:H:186:LYS:HE2	1:H:204:GLU:CD	2.26	0.55
1:A:152:VAL:HG13	1:A:165:ILE:HD12	1.89	0.54
1:D:43:TYR:CE1	1:D:147:LEU:HD11	2.42	0.54
1:H:69:TRP:C	1:H:71:TYR:H	2.11	0.54
1:F:178:ARG:HG3	1:F:182:TRP:CE2	2.42	0.54
1:F:142:GLY:C	1:F:145:PRO:HD2	2.29	0.53
1:H:14:LEU:HD11	1:H:84:ALA:HA	1.90	0.53
1:E:137:SER:O	1:E:141:ARG:HG3	2.07	0.53
1:F:30:ARG:HD2	1:F:30:ARG:C	2.29	0.53
1:H:186:LYS:HE2	1:H:204:GLU:OE2	2.09	0.53
1:E:141:ARG:HG2	1:E:141:ARG:HH11	1.73	0.53
1:D:116:ILE:HD11	1:D:200:LEU:HD11	1.91	0.53
1:H:60:MET:HG3	1:H:78:TRP:HZ2	1.71	0.53
1:G:190:GLU:OE1	1:G:197:ARG:NH1	2.42	0.53
1:A:51:THR:HG21	1:D:119:TYR:CE2	2.44	0.53
1:A:142:GLY:C	1:A:145:PRO:HD2	2.29	0.52
1:C:192:ASP:CG	1:C:197:ARG:HD2	2.30	0.52
1:G:81:HIS:HE1	1:G:107:GLU:OE2	1.92	0.52
1:B:164:ARG:HG3	1:B:202:ALA:HB1	1.90	0.52
1:A:171:HIS:CD2	1:A:197:ARG:HB3	2.44	0.52
1:H:21:ARG:HG3	3:H:241:HOH:O	2.09	0.52
1:B:77:ARG:HA	1:B:80:GLN:HE21	1.74	0.52
1:A:30:ARG:HB2	1:A:30:ARG:HH11	1.74	0.52
1:H:144:GLY:O	1:H:148:LEU:HD23	2.10	0.52
1:F:37:PRO:HG2	1:F:45:LEU:HD13	1.92	0.52
1:H:127:LEU:O	1:H:165:ILE:HA	2.09	0.51
1:C:136:LEU:CD1	1:C:136:LEU:H	2.23	0.51
1:A:186:LYS:HE3	1:A:188:LEU:HD11	1.91	0.51
1:E:12:ASP:O	1:E:13:ALA:HB2	2.09	0.51
1:A:41:PRO:HG3	1:D:186:LYS:HE3	1.92	0.51
1:A:10:ALA:HA	1:E:77:ARG:NH1	2.23	0.51
1:D:81:HIS:HD2	3:D:741:HOH:O	1.92	0.51
1:G:51:THR:HG23	3:G:758:HOH:O	2.10	0.51
1:B:115:LEU:CD1	1:B:194:THR:HG23	2.41	0.50
1:C:54:GLU:N	1:C:54:GLU:OE2	2.43	0.50
1:D:193:THR:HG22	1:D:196:ARG:O	2.11	0.50
1:D:137:SER:O	1:D:141:ARG:HG2	2.11	0.50
1:A:30:ARG:HB2	1:A:30:ARG:NH1	2.26	0.50
1:B:69:TRP:O	1:B:70:GLU:HB2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:MET:HG3	1:B:78:TRP:HZ2	1.70	0.49
1:B:60:MET:HE2	1:B:65:LEU:HD13	1.94	0.49
1:C:136:LEU:N	1:C:136:LEU:HD12	2.24	0.49
1:C:73:TRP:CE3	1:C:77:ARG:HD2	2.47	0.49
1:C:77:ARG:NH1	1:C:77:ARG:HG3	2.27	0.49
1:G:173:ASN:O	1:G:177:ARG:HG3	2.13	0.49
1:A:160:PRO:HD2	1:B:67:ALA:CB	2.43	0.49
1:G:49:GLN:NE2	1:G:49:GLN:HA	2.27	0.49
1:A:123:ASP:OD1	1:A:124:PRO:HD2	2.13	0.49
1:E:115:LEU:HD23	1:E:116:ILE:H	1.78	0.49
1:F:178:ARG:HG3	1:F:182:TRP:CZ2	2.47	0.48
1:B:69:TRP:NE1	1:B:132:ALA:HB2	2.28	0.48
1:E:144:GLY:N	1:E:145:PRO:HD2	2.28	0.48
1:D:172:ARG:HH11	1:D:172:ARG:CB	2.26	0.48
1:B:100:GLY:HA2	1:C:187:PHE:HB2	1.95	0.48
1:C:140:ASN:HB3	1:C:141:ARG:NH2	2.28	0.48
1:D:103:GLY:HA3	2:D:701:BOG:H2'2	1.95	0.48
1:C:60:MET:HG3	1:C:78:TRP:HZ2	1.71	0.48
1:H:61:ASN:C	1:H:62:ARG:HH11	2.17	0.48
1:G:127:LEU:O	1:G:165:ILE:HA	2.14	0.48
1:E:43:TYR:CE1	1:E:147:LEU:HD11	2.49	0.48
1:D:175:ALA:O	1:D:179:LEU:HG	2.14	0.47
1:H:146:LEU:HG	3:H:272:HOH:O	2.12	0.47
1:D:114:ASP:HB3	1:D:116:ILE:HG22	1.96	0.47
1:E:39:LEU:HD21	1:E:151:ILE:HG12	1.95	0.47
1:F:77:ARG:CG	1:F:77:ARG:NH1	2.75	0.47
1:B:73:TRP:HB3	1:B:77:ARG:HB3	1.95	0.47
1:E:144:GLY:O	1:E:148:LEU:HD13	2.15	0.47
1:C:14:LEU:HD11	1:C:84:ALA:HA	1.95	0.47
1:B:55:MET:O	1:B:58:GLU:HB3	2.13	0.47
1:B:193:THR:O	1:B:194:THR:C	2.53	0.47
1:H:15:VAL:HG22	3:H:253:HOH:O	2.15	0.47
1:B:115:LEU:HD13	3:B:262:HOH:O	2.14	0.47
1:C:29:ARG:HD3	3:C:257:HOH:O	2.14	0.47
1:E:190:GLU:HB3	1:E:197:ARG:NH1	2.30	0.47
1:H:60:MET:HE1	1:H:65:LEU:HD13	1.97	0.47
1:B:65:LEU:HD22	1:B:132:ALA:HB1	1.97	0.47
1:F:60:MET:HG3	1:F:78:TRP:HZ2	1.78	0.46
1:G:171:HIS:NE2	1:G:172:ARG:HG3	2.30	0.46
1:F:188:LEU:HD12	1:F:188:LEU:N	2.29	0.46
1:E:186:LYS:HZ1	1:E:204:GLU:HG3	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:30:ARG:HH11	1:C:30:ARG:HG3	1.79	0.46
1:A:54:GLU:CD	1:A:54:GLU:H	2.18	0.46
1:F:8:GLY:N	1:F:79:ARG:HD3	2.31	0.46
1:B:69:TRP:HE1	1:B:132:ALA:HB2	1.80	0.46
1:D:116:ILE:HD11	1:D:200:LEU:HD13	1.95	0.46
1:E:178:ARG:HD3	1:E:178:ARG:HA	1.75	0.46
1:F:114:ASP:OD1	1:F:116:ILE:HG23	2.16	0.46
1:A:194:THR:HG23	1:A:195:ASN:N	2.31	0.46
1:E:120:TYR:CE1	1:E:122:ALA:HA	2.50	0.46
1:A:78:TRP:O	1:A:82:LEU:HG	2.15	0.45
1:A:40:GLU:OE1	1:A:40:GLU:HA	2.15	0.45
1:E:109:TYR:CZ	1:E:128:GLY:HA3	2.51	0.45
1:D:188:LEU:O	1:D:191:HIS:HE1	1.98	0.45
1:H:119:TYR:CZ	1:H:193:THR:HG22	2.51	0.45
1:A:174:THR:HG22	1:A:178:ARG:NE	2.31	0.45
1:C:50:LEU:HB3	1:C:79:ARG:NH1	2.30	0.45
1:E:30:ARG:NH1	1:E:30:ARG:HG3	2.30	0.45
1:C:73:TRP:HB3	1:C:77:ARG:CB	2.46	0.45
1:H:62:ARG:HB3	1:H:63:PRO:HD2	1.99	0.45
1:A:51:THR:OG1	1:D:191:HIS:HD2	2.00	0.45
1:B:186:LYS:HE2	3:B:243:HOH:O	2.16	0.45
1:A:73:TRP:CE3	1:A:77:ARG:HG3	2.52	0.45
1:D:152:VAL:HG11	1:D:203:LEU:HD23	1.99	0.45
1:H:62:ARG:HD3	1:H:62:ARG:HA	1.81	0.45
1:B:63:PRO:HB2	3:B:256:HOH:O	2.16	0.45
1:H:34:PRO:O	1:H:92:LEU:HD22	2.17	0.45
1:C:105:TYR:C	1:C:106:LEU:HD12	2.37	0.45
1:D:178:ARG:HD2	1:D:182:TRP:CH2	2.51	0.44
1:E:106:LEU:HD11	1:E:151:ILE:HD13	1.98	0.44
1:C:194:THR:O	1:C:195:ASN:CB	2.65	0.44
1:A:47:VAL:HG23	3:A:233:HOH:O	2.17	0.44
1:D:205:ALA:HB1	1:D:206:PRO:HD2	2.00	0.44
1:A:127:LEU:O	1:A:165:ILE:HA	2.17	0.44
1:A:136:LEU:O	1:A:136:LEU:HD23	2.16	0.44
1:D:185:CYS:HB3	1:D:201:TYR:HB3	2.00	0.44
1:E:171:HIS:HB3	1:E:197:ARG:HG2	1.98	0.44
1:A:61:ASN:OD1	1:A:72:ASP:HB2	2.17	0.44
1:B:56:LEU:HD13	1:B:78:TRP:CZ3	2.53	0.44
1:A:65:LEU:HD22	1:A:132:ALA:HB1	1.98	0.44
1:C:109:TYR:CZ	1:C:128:GLY:HA3	2.53	0.44
1:G:49:GLN:HE21	1:G:49:GLN:HA	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:123:ASP:HB3	1:G:126:ASP:OD1	2.17	0.43
1:B:124:PRO:HB2	1:B:125:TYR:CE1	2.53	0.43
1:F:145:PRO:HD3	3:F:280:HOH:O	2.18	0.43
1:E:143:PHE:CZ	2:E:702:BOG:H4'2	2.54	0.43
1:H:156:PHE:CD1	1:H:206:PRO:HD3	2.54	0.43
1:C:186:LYS:HE2	1:C:204:GLU:HG3	2.00	0.43
1:F:187:PHE:CZ	1:F:189:GLY:HA2	2.54	0.43
1:C:208:THR:O	1:C:209:ALA:HB3	2.19	0.43
1:A:106:LEU:HD21	1:A:151:ILE:HD13	2.01	0.43
1:D:190:GLU:HA	1:D:198:MET:O	2.18	0.43
1:B:71:TYR:O	1:B:73:TRP:N	2.48	0.43
1:G:196:ARG:HH11	1:G:196:ARG:CG	2.32	0.43
1:C:49:GLN:O	1:C:52:ASP:HB2	2.18	0.43
1:A:24:LEU:HA	1:A:25:PRO:HD3	1.88	0.43
1:E:143:PHE:CE1	2:E:702:BOG:H4'2	2.54	0.43
1:B:40:GLU:OE1	1:B:40:GLU:HA	2.19	0.43
1:A:173:ASN:O	1:A:177:ARG:HG3	2.18	0.43
1:F:190:GLU:HA	1:F:198:MET:O	2.19	0.43
1:E:27:GLN:HG3	1:E:30:ARG:NH2	2.33	0.43
1:H:11:ASP:CG	1:H:79:ARG:HH12	2.21	0.42
1:A:85:GLN:NE2	1:A:91:SER:OG	2.47	0.42
1:G:49:GLN:NE2	3:G:783:HOH:O	2.51	0.42
1:F:173:ASN:O	1:F:177:ARG:HG3	2.19	0.42
1:B:125:TYR:CD1	1:B:125:TYR:N	2.87	0.42
1:G:101:THR:HG21	2:G:703:BOG:H2	2.00	0.42
1:A:85:GLN:HE22	1:A:107:GLU:HG2	1.83	0.42
1:E:185:CYS:HB3	1:E:201:TYR:HB3	2.01	0.42
1:B:137:SER:O	1:B:141:ARG:HB2	2.19	0.42
1:A:137:SER:HA	1:A:141:ARG:NH1	2.35	0.42
1:A:85:GLN:HG2	1:A:91:SER:HB3	2.01	0.42
1:B:178:ARG:HG2	3:B:263:HOH:O	2.19	0.42
1:H:69:TRP:C	1:H:71:TYR:N	2.73	0.42
1:E:31:LEU:HD21	1:E:125:TYR:CE2	2.54	0.42
1:G:30:ARG:HH11	1:G:30:ARG:HG2	1.85	0.42
1:A:65:LEU:CD2	1:A:132:ALA:HB1	2.50	0.42
1:G:35:PRO:HB2	1:G:158:ASN:HB3	2.01	0.42
1:G:178:ARG:HH11	1:G:178:ARG:HG3	1.85	0.42
1:A:148:LEU:HB2	1:A:149:PRO:HD3	2.02	0.42
1:H:45:LEU:HD23	1:H:96:GLY:HA2	2.02	0.42
1:C:145:PRO:HB3	1:C:179:LEU:HD21	2.02	0.42
1:B:100:GLY:HA2	1:C:187:PHE:CB	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:104:GLY:HA3	2:E:702:BOG:H7'1	2.02	0.42
1:A:60:MET:HG2	1:A:78:TRP:CZ2	2.46	0.41
1:F:14:LEU:HD11	1:F:84:ALA:HA	2.02	0.41
1:B:15:VAL:HG13	3:B:212:HOH:O	2.19	0.41
1:B:186:LYS:HD2	1:B:188:LEU:CD2	2.50	0.41
1:F:73:TRP:HB3	1:F:74:PRO:HD2	2.01	0.41
1:B:193:THR:HG22	1:B:196:ARG:O	2.20	0.41
1:H:142:GLY:O	1:H:145:PRO:HD2	2.21	0.41
1:C:106:LEU:HD23	1:C:129:LEU:HD11	2.03	0.41
1:D:169:PRO:HD2	1:D:199:ALA:O	2.21	0.41
1:H:64:HIS:N	1:H:64:HIS:CD2	2.88	0.41
1:B:203:LEU:C	1:B:203:LEU:HD23	2.41	0.41
1:C:73:TRP:CB	1:C:77:ARG:HB3	2.49	0.41
1:B:24:LEU:HD13	1:B:113:LYS:HG2	2.01	0.41
1:A:79:ARG:HH11	1:A:79:ARG:HG2	1.85	0.41
1:E:172:ARG:HD2	1:H:163:ARG:NE	2.36	0.41
1:D:177:ARG:O	1:D:181:GLU:HG3	2.21	0.41
1:H:73:TRP:HB3	1:H:77:ARG:HB3	2.03	0.41
1:E:186:LYS:HG3	1:E:204:GLU:OE2	2.21	0.41
1:A:34:PRO:HA	1:A:35:PRO:HD3	1.84	0.41
1:G:188:LEU:O	1:G:191:HIS:HE1	2.03	0.41
1:C:73:TRP:CD2	1:C:77:ARG:HD2	2.56	0.41
1:D:65:LEU:HD23	1:D:65:LEU:HA	1.94	0.41
1:E:159:GLU:OE1	1:E:161:ARG:NH2	2.53	0.41
1:E:40:GLU:OE2	1:E:40:GLU:HA	2.20	0.41
1:H:73:TRP:HB3	1:H:74:PRO:CD	2.49	0.41
1:G:171:HIS:CD2	1:G:172:ARG:HG3	2.56	0.41
1:A:34:PRO:O	1:A:92:LEU:HD22	2.21	0.41
1:E:142:GLY:O	1:E:145:PRO:HG2	2.22	0.40
1:E:190:GLU:HB3	1:E:197:ARG:HH11	1.86	0.40
1:G:65:LEU:HA	1:G:65:LEU:HD23	1.86	0.40
1:C:66:ALA:O	1:C:70:GLU:HA	2.21	0.40
1:F:187:PHE:C	1:F:188:LEU:HD12	2.41	0.40
1:D:109:TYR:CZ	1:D:128:GLY:HA3	2.57	0.40
1:C:171:HIS:CD2	1:C:172:ARG:HG3	2.56	0.40
1:C:190:GLU:HA	1:C:198:MET:O	2.22	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	196/210 (93%)	187 (95%)	6 (3%)	3 (2%)	13	9
1	B	189/210 (90%)	183 (97%)	4 (2%)	2 (1%)	17	14
1	C	197/210 (94%)	188 (95%)	7 (4%)	2 (1%)	19	16
1	D	198/210 (94%)	192 (97%)	3 (2%)	3 (2%)	13	9
1	E	194/210 (92%)	184 (95%)	8 (4%)	2 (1%)	19	16
1	F	200/210 (95%)	190 (95%)	9 (4%)	1 (0%)	34	35
1	G	199/210 (95%)	194 (98%)	4 (2%)	1 (0%)	34	35
1	H	195/210 (93%)	190 (97%)	3 (2%)	2 (1%)	19	16
All	All	1568/1680 (93%)	1508 (96%)	44 (3%)	16 (1%)	19	16

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	194	THR
1	E	13	ALA
1	F	208	THR
1	A	11	ASP
1	D	195	ASN
1	H	70	GLU
1	H	195	ASN
1	A	194	THR
1	G	194	THR
1	B	72	ASP
1	C	12	ASP
1	D	208	THR
1	D	194	THR
1	E	206	PRO
1	A	63	PRO
1	C	63	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	162/170 (95%)	155 (96%)	7 (4%)	35	43
1	B	158/170 (93%)	149 (94%)	9 (6%)	25	29
1	C	163/170 (96%)	155 (95%)	8 (5%)	31	36
1	D	163/170 (96%)	159 (98%)	4 (2%)	55	67
1	E	161/170 (95%)	159 (99%)	2 (1%)	78	88
1	F	164/170 (96%)	159 (97%)	5 (3%)	48	60
1	G	163/170 (96%)	159 (98%)	4 (2%)	55	67
1	H	161/170 (95%)	153 (95%)	8 (5%)	30	35
All	All	1295/1360 (95%)	1248 (96%)	47 (4%)	42	52

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	ASP
1	A	30	ARG
1	A	64	HIS
1	A	188	LEU
1	A	197	ARG
1	A	198	MET
1	A	203	LEU
1	B	40	GLU
1	B	63	PRO
1	B	90	TYR
1	B	115	LEU
1	B	121	ASP
1	B	136	LEU
1	B	143	PHE
1	B	164	ARG
1	B	198	MET
1	C	26	ASP
1	C	77	ARG
1	C	90	TYR

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Mol	Chain	Res	Type
1	C	135	ASP
1	C	141	ARG
1	C	161	ARG
1	C	171	HIS
1	C	197	ARG
1	D	54	GLU
1	D	90	TYR
1	D	192	ASP
1	D	195	ASN
1	E	90	TYR
1	E	121	ASP
1	F	51	THR
1	F	77	ARG
1	F	90	TYR
1	F	148	LEU
1	F	186	LYS
1	G	90	TYR
1	G	195	ASN
1	G	198	MET
1	G	207	THR
1	H	24	LEU
1	H	33	ARG
1	H	51	THR
1	H	54	GLU
1	H	64	HIS
1	H	90	TYR
1	H	101	THR
1	H	135	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	HIS
1	A	81	HIS
1	A	85	GLN
1	A	130	HIS
1	A	140	ASN
1	A	171	HIS
1	B	80	GLN
1	B	83	ASN
1	B	85	GLN
1	B	99	HIS

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Mol	Chain	Res	Type
1	B	118	HIS
1	B	158	ASN
1	B	171	HIS
1	C	27	GLN
1	C	85	GLN
1	C	118	HIS
1	C	140	ASN
1	C	158	ASN
1	C	195	ASN
1	D	64	HIS
1	D	81	HIS
1	D	85	GLN
1	D	99	HIS
1	D	118	HIS
1	D	171	HIS
1	D	191	HIS
1	E	85	GLN
1	E	99	HIS
1	E	158	ASN
1	F	27	GLN
1	F	85	GLN
1	F	118	HIS
1	F	140	ASN
1	F	158	ASN
1	F	171	HIS
1	F	191	HIS
1	G	49	GLN
1	G	64	HIS
1	G	81	HIS
1	G	85	GLN
1	G	191	HIS
1	G	195	ASN
1	H	64	HIS
1	H	85	GLN
1	H	118	HIS
1	H	130	HIS
1	H	171	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

3 ligands 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	BOG	D	701	-	20,20,20	1.16	3 (15%)	25,25,25	0.63	0
2	BOG	E	702	-	20,20,20	1.19	2 (10%)	25,25,25	0.66	0
2	BOG	G	703	-	20,20,20	1.06	2 (10%)	25,25,25	0.66	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
2	BOG	D	701	-	-	0/11/31/31	0/1/1/1
2	BOG	E	702	-	-	0/11/31/31	0/1/1/1
2	BOG	G	703	-	-	0/11/31/31	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	701	BOG	O1-C1	2.01	1.43	1.40
2	G	703	BOG	C4-C5	2.05	1.57	1.53
2	D	701	BOG	C4-C5	2.30	1.58	1.53
2	E	702	BOG	C4-C5	2.43	1.58	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	703	BOG	O5-C1	2.90	1.49	1.41
2	D	701	BOG	O5-C1	2.99	1.49	1.41
2	E	702	BOG	O5-C1	3.04	1.49	1.41

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	701	BOG	3	0
2	E	702	BOG	4	0
2	G	703	BOG	2	0

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	198/210 (94%)	0.35	16 (8%) 15 14	23, 37, 76, 82	0
1	B	193/210 (91%)	0.34	12 (6%) 24 23	21, 38, 61, 77	0
1	C	199/210 (94%)	0.17	12 (6%) 25 25	16, 31, 65, 74	0
1	D	200/210 (95%)	0.19	11 (5%) 29 28	18, 32, 59, 79	0
1	E	196/210 (93%)	0.40	16 (8%) 14 14	19, 36, 61, 76	0
1	F	202/210 (96%)	0.08	9 (4%) 37 36	12, 24, 48, 84	0
1	G	201/210 (95%)	-0.03	8 (3%) 42 41	13, 25, 43, 71	0
1	H	197/210 (93%)	0.35	22 (11%) 7 6	24, 39, 81, 90	0
All	All	1586/1680 (94%)	0.23	106 (6%) 21 20	12, 32, 65, 90	0

All (106) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	12	ASP	8.9
1	F	8	GLY	8.9
1	C	209	ALA	8.4
1	D	208	THR	8.2
1	A	10	ALA	8.1
1	F	10	ALA	7.8
1	F	9	GLN	6.4
1	E	206	PRO	6.3
1	A	11	ASP	6.1
1	D	10	ALA	6.1
1	C	208	THR	5.8
1	E	207	THR	5.4
1	F	209	ALA	5.4
1	D	11	ASP	5.3
1	A	136	LEU	5.2
1	A	12	ASP	4.9

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Mol	Chain	Res	Type	RSRZ
1	B	13	ALA	4.7
1	B	207	THR	4.6
1	H	10	ALA	4.5
1	F	208	THR	4.4
1	H	69	TRP	4.3
1	G	11	ASP	4.2
1	B	194	THR	4.2
1	G	10	ALA	4.1
1	B	143	PHE	4.1
1	A	68	ALA	4.0
1	F	11	ASP	4.0
1	F	194	THR	3.9
1	C	140	ASN	3.8
1	A	137	SER	3.8
1	E	140	ASN	3.8
1	H	11	ASP	3.8
1	B	99	HIS	3.7
1	H	140	ASN	3.7
1	D	207	THR	3.7
1	A	66	ALA	3.6
1	C	207	THR	3.6
1	B	135	ASP	3.5
1	A	71	TYR	3.5
1	A	75	ALA	3.5
1	A	194	THR	3.4
1	B	12	ASP	3.4
1	C	12	ASP	3.4
1	E	30	ARG	3.3
1	H	141	ARG	3.3
1	A	69	TRP	3.3
1	C	141	ARG	3.1
1	H	194	THR	3.1
1	E	12	ASP	3.1
1	D	209	ALA	3.0
1	D	30	ARG	3.0
1	B	26	ASP	3.0
1	E	26	ASP	3.0
1	B	137	SER	2.9
1	G	194	THR	2.9
1	B	136	LEU	2.9
1	D	178	ARG	2.9
1	E	163	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	11	ASP	2.9
1	H	65	LEU	2.9
1	C	66	ALA	2.8
1	A	140	ASN	2.8
1	C	75	ALA	2.8
1	D	26	ASP	2.7
1	D	163	ARG	2.6
1	H	98	TRP	2.6
1	C	194	THR	2.6
1	A	135	ASP	2.6
1	E	13	ALA	2.6
1	G	13	ALA	2.6
1	D	140	ASN	2.5
1	H	12	ASP	2.5
1	F	195	ASN	2.5
1	H	63	PRO	2.5
1	H	135	ASP	2.5
1	E	157	ALA	2.5
1	H	71	TYR	2.5
1	A	63	PRO	2.5
1	G	210	ALA	2.4
1	E	125	TYR	2.4
1	G	26	ASP	2.4
1	H	138	LYS	2.4
1	A	99	HIS	2.4
1	H	99	HIS	2.4
1	E	174	THR	2.4
1	B	63	PRO	2.3
1	H	131	ALA	2.3
1	H	30	ARG	2.3
1	D	29	ARG	2.3
1	A	13	ALA	2.2
1	C	61	ASN	2.2
1	F	129	LEU	2.2
1	E	131	ALA	2.2
1	C	73	TRP	2.2
1	G	208	THR	2.2
1	E	195	ASN	2.1
1	H	143	PHE	2.1
1	H	13	ALA	2.1
1	H	29	ARG	2.1
1	H	62	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	205	ALA	2.1
1	H	195	ASN	2.1
1	E	172	ARG	2.1
1	B	30	ARG	2.1
1	E	156	PHE	2.0
1	H	206	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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(\AA^2)	Q<0.9
2	BOG	D	701	20/20	0.70	0.29	5.63	51,60,63,66	0
2	BOG	G	703	20/20	0.71	0.30	4.63	50,58,61,62	0
2	BOG	E	702	20/20	0.74	0.27	3.43	52,61,63,63	0

6.5 Other polymers [i](#)

There are no such residues in this entry.