



Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 12:21 AM GMT

PDB ID : 3KTM
Title : Structure of the Heparin-induced E1-Dimer of the Amyloid Precursor Protein (APP)
Authors : Dahms, S.O.; Hoefgen, S.; Roeser, D.; Schlott, B.; Guhrs, K.H.; Than, M.E.
Deposited on : 2009-11-25
Resolution : 2.70 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>

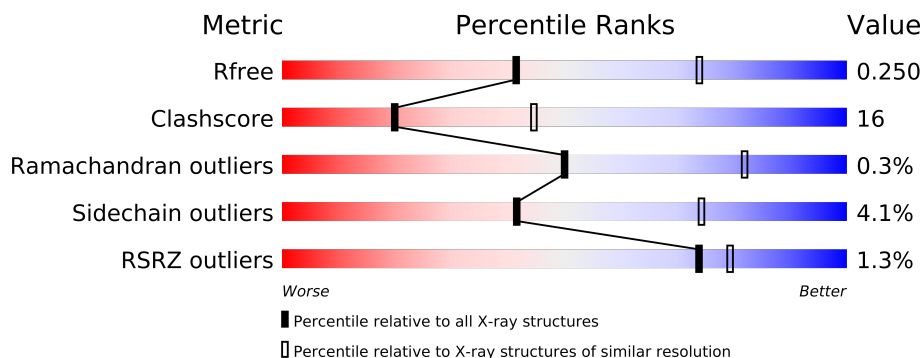
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.70 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	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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	191	
1	B	191	
1	C	191	
1	D	191	
1	E	191	
1	F	191	
1	G	191	
1	H	191	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	BU4	D	2	-	X

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Geometry	Electron density
2	BU4	F	4	-	X
2	BU4	G	3	-	X
4	ACT	B	208	-	X
4	ACT	C	8	-	X
4	ACT	E	9	-	X
4	ACT	H	6	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11436 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 Amyloid beta A4 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	175	Total	C	N	O	S	32	0	0
			1392	878	245	252	17			
1	B	173	Total	C	N	O	S	27	0	0
			1369	863	241	248	17			
1	C	174	Total	C	N	O	S	20	0	0
			1374	866	242	249	17			
1	D	173	Total	C	N	O	S	26	0	0
			1373	867	241	248	17			
1	E	177	Total	C	N	O	S	35	0	0
			1401	883	247	254	17			
1	F	175	Total	C	N	O	S	37	0	0
			1392	878	245	252	17			
1	G	174	Total	C	N	O	S	50	0	0
			1382	872	242	251	17			
1	H	175	Total	C	N	O	S	38	0	0
			1382	872	243	250	17			

There are 144 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	MET	-	INITIATING METHIONINE	UNP P05067
A	191	ILE	-	EXPRESSION TAG	UNP P05067
A	192	GLU	-	EXPRESSION TAG	UNP P05067
A	193	GLY	-	EXPRESSION TAG	UNP P05067
A	194	ARG	-	EXPRESSION TAG	UNP P05067
A	195	LYS	-	EXPRESSION TAG	UNP P05067
A	196	LEU	-	EXPRESSION TAG	UNP P05067
A	197	ALA	-	EXPRESSION TAG	UNP P05067
A	198	ALA	-	EXPRESSION TAG	UNP P05067
A	199	ALA	-	EXPRESSION TAG	UNP P05067
A	200	LEU	-	EXPRESSION TAG	UNP P05067
A	201	GLU	-	EXPRESSION TAG	UNP P05067
A	202	HIS	-	EXPRESSION TAG	UNP P05067

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	203	HIS	-	EXPRESSION TAG	UNP P05067
A	204	HIS	-	EXPRESSION TAG	UNP P05067
A	205	HIS	-	EXPRESSION TAG	UNP P05067
A	206	HIS	-	EXPRESSION TAG	UNP P05067
A	207	HIS	-	EXPRESSION TAG	UNP P05067
B	17	MET	-	INITIATING METHIONINE	UNP P05067
B	191	ILE	-	EXPRESSION TAG	UNP P05067
B	192	GLU	-	EXPRESSION TAG	UNP P05067
B	193	GLY	-	EXPRESSION TAG	UNP P05067
B	194	ARG	-	EXPRESSION TAG	UNP P05067
B	195	LYS	-	EXPRESSION TAG	UNP P05067
B	196	LEU	-	EXPRESSION TAG	UNP P05067
B	197	ALA	-	EXPRESSION TAG	UNP P05067
B	198	ALA	-	EXPRESSION TAG	UNP P05067
B	199	ALA	-	EXPRESSION TAG	UNP P05067
B	200	LEU	-	EXPRESSION TAG	UNP P05067
B	201	GLU	-	EXPRESSION TAG	UNP P05067
B	202	HIS	-	EXPRESSION TAG	UNP P05067
B	203	HIS	-	EXPRESSION TAG	UNP P05067
B	204	HIS	-	EXPRESSION TAG	UNP P05067
B	205	HIS	-	EXPRESSION TAG	UNP P05067
B	206	HIS	-	EXPRESSION TAG	UNP P05067
B	207	HIS	-	EXPRESSION TAG	UNP P05067
C	17	MET	-	INITIATING METHIONINE	UNP P05067
C	191	ILE	-	EXPRESSION TAG	UNP P05067
C	192	GLU	-	EXPRESSION TAG	UNP P05067
C	193	GLY	-	EXPRESSION TAG	UNP P05067
C	194	ARG	-	EXPRESSION TAG	UNP P05067
C	195	LYS	-	EXPRESSION TAG	UNP P05067
C	196	LEU	-	EXPRESSION TAG	UNP P05067
C	197	ALA	-	EXPRESSION TAG	UNP P05067
C	198	ALA	-	EXPRESSION TAG	UNP P05067
C	199	ALA	-	EXPRESSION TAG	UNP P05067
C	200	LEU	-	EXPRESSION TAG	UNP P05067
C	201	GLU	-	EXPRESSION TAG	UNP P05067
C	202	HIS	-	EXPRESSION TAG	UNP P05067
C	203	HIS	-	EXPRESSION TAG	UNP P05067
C	204	HIS	-	EXPRESSION TAG	UNP P05067
C	205	HIS	-	EXPRESSION TAG	UNP P05067
C	206	HIS	-	EXPRESSION TAG	UNP P05067
C	207	HIS	-	EXPRESSION TAG	UNP P05067
D	17	MET	-	INITIATING METHIONINE	UNP P05067

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	191	ILE	-	EXPRESSION TAG	UNP P05067
D	192	GLU	-	EXPRESSION TAG	UNP P05067
D	193	GLY	-	EXPRESSION TAG	UNP P05067
D	194	ARG	-	EXPRESSION TAG	UNP P05067
D	195	LYS	-	EXPRESSION TAG	UNP P05067
D	196	LEU	-	EXPRESSION TAG	UNP P05067
D	197	ALA	-	EXPRESSION TAG	UNP P05067
D	198	ALA	-	EXPRESSION TAG	UNP P05067
D	199	ALA	-	EXPRESSION TAG	UNP P05067
D	200	LEU	-	EXPRESSION TAG	UNP P05067
D	201	GLU	-	EXPRESSION TAG	UNP P05067
D	202	HIS	-	EXPRESSION TAG	UNP P05067
D	203	HIS	-	EXPRESSION TAG	UNP P05067
D	204	HIS	-	EXPRESSION TAG	UNP P05067
D	205	HIS	-	EXPRESSION TAG	UNP P05067
D	206	HIS	-	EXPRESSION TAG	UNP P05067
D	207	HIS	-	EXPRESSION TAG	UNP P05067
E	17	MET	-	INITIATING METHIONINE	UNP P05067
E	191	ILE	-	EXPRESSION TAG	UNP P05067
E	192	GLU	-	EXPRESSION TAG	UNP P05067
E	193	GLY	-	EXPRESSION TAG	UNP P05067
E	194	ARG	-	EXPRESSION TAG	UNP P05067
E	195	LYS	-	EXPRESSION TAG	UNP P05067
E	196	LEU	-	EXPRESSION TAG	UNP P05067
E	197	ALA	-	EXPRESSION TAG	UNP P05067
E	198	ALA	-	EXPRESSION TAG	UNP P05067
E	199	ALA	-	EXPRESSION TAG	UNP P05067
E	200	LEU	-	EXPRESSION TAG	UNP P05067
E	201	GLU	-	EXPRESSION TAG	UNP P05067
E	202	HIS	-	EXPRESSION TAG	UNP P05067
E	203	HIS	-	EXPRESSION TAG	UNP P05067
E	204	HIS	-	EXPRESSION TAG	UNP P05067
E	205	HIS	-	EXPRESSION TAG	UNP P05067
E	206	HIS	-	EXPRESSION TAG	UNP P05067
E	207	HIS	-	EXPRESSION TAG	UNP P05067
F	17	MET	-	INITIATING METHIONINE	UNP P05067
F	191	ILE	-	EXPRESSION TAG	UNP P05067
F	192	GLU	-	EXPRESSION TAG	UNP P05067
F	193	GLY	-	EXPRESSION TAG	UNP P05067
F	194	ARG	-	EXPRESSION TAG	UNP P05067
F	195	LYS	-	EXPRESSION TAG	UNP P05067
F	196	LEU	-	EXPRESSION TAG	UNP P05067

Continued on next page...

Continued from previous page...

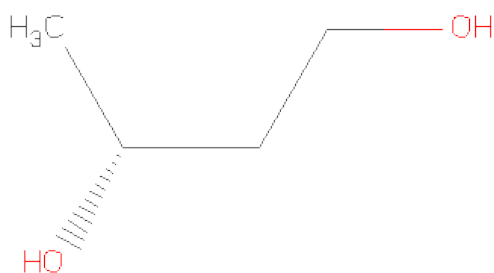
Chain	Residue	Modelled	Actual	Comment	Reference
F	197	ALA	-	EXPRESSION TAG	UNP P05067
F	198	ALA	-	EXPRESSION TAG	UNP P05067
F	199	ALA	-	EXPRESSION TAG	UNP P05067
F	200	LEU	-	EXPRESSION TAG	UNP P05067
F	201	GLU	-	EXPRESSION TAG	UNP P05067
F	202	HIS	-	EXPRESSION TAG	UNP P05067
F	203	HIS	-	EXPRESSION TAG	UNP P05067
F	204	HIS	-	EXPRESSION TAG	UNP P05067
F	205	HIS	-	EXPRESSION TAG	UNP P05067
F	206	HIS	-	EXPRESSION TAG	UNP P05067
F	207	HIS	-	EXPRESSION TAG	UNP P05067
G	17	MET	-	INITIATING METHIONINE	UNP P05067
G	191	ILE	-	EXPRESSION TAG	UNP P05067
G	192	GLU	-	EXPRESSION TAG	UNP P05067
G	193	GLY	-	EXPRESSION TAG	UNP P05067
G	194	ARG	-	EXPRESSION TAG	UNP P05067
G	195	LYS	-	EXPRESSION TAG	UNP P05067
G	196	LEU	-	EXPRESSION TAG	UNP P05067
G	197	ALA	-	EXPRESSION TAG	UNP P05067
G	198	ALA	-	EXPRESSION TAG	UNP P05067
G	199	ALA	-	EXPRESSION TAG	UNP P05067
G	200	LEU	-	EXPRESSION TAG	UNP P05067
G	201	GLU	-	EXPRESSION TAG	UNP P05067
G	202	HIS	-	EXPRESSION TAG	UNP P05067
G	203	HIS	-	EXPRESSION TAG	UNP P05067
G	204	HIS	-	EXPRESSION TAG	UNP P05067
G	205	HIS	-	EXPRESSION TAG	UNP P05067
G	206	HIS	-	EXPRESSION TAG	UNP P05067
G	207	HIS	-	EXPRESSION TAG	UNP P05067
H	17	MET	-	INITIATING METHIONINE	UNP P05067
H	191	ILE	-	EXPRESSION TAG	UNP P05067
H	192	GLU	-	EXPRESSION TAG	UNP P05067
H	193	GLY	-	EXPRESSION TAG	UNP P05067
H	194	ARG	-	EXPRESSION TAG	UNP P05067
H	195	LYS	-	EXPRESSION TAG	UNP P05067
H	196	LEU	-	EXPRESSION TAG	UNP P05067
H	197	ALA	-	EXPRESSION TAG	UNP P05067
H	198	ALA	-	EXPRESSION TAG	UNP P05067
H	199	ALA	-	EXPRESSION TAG	UNP P05067
H	200	LEU	-	EXPRESSION TAG	UNP P05067
H	201	GLU	-	EXPRESSION TAG	UNP P05067
H	202	HIS	-	EXPRESSION TAG	UNP P05067

Continued on next page...

Continued from previous page...

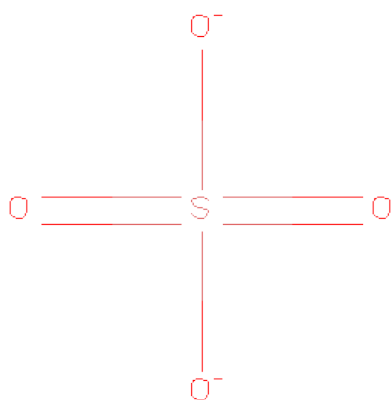
Chain	Residue	Modelled	Actual	Comment	Reference
H	203	HIS	-	EXPRESSION TAG	UNP P05067
H	204	HIS	-	EXPRESSION TAG	UNP P05067
H	205	HIS	-	EXPRESSION TAG	UNP P05067
H	206	HIS	-	EXPRESSION TAG	UNP P05067
H	207	HIS	-	EXPRESSION TAG	UNP P05067

- Molecule 2 is (3R)-BUTANE-1,3-DIOL (three-letter code: BU4) (formula: C₄H₁₀O₂).



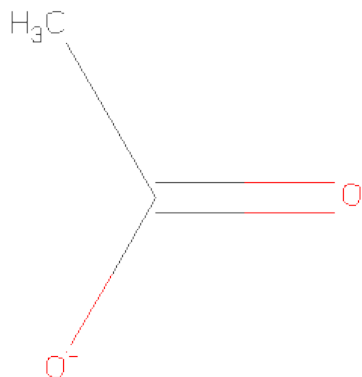
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	4	2		
2	B	1	Total	C	O	0	0
			6	4	2		
2	C	1	Total	C	O	0	0
			6	4	2		
2	D	1	Total	C	O	0	0
			6	4	2		
2	E	1	Total	C	O	0	0
			6	4	2		
2	F	1	Total	C	O	0	0
			6	4	2		
2	G	1	Total	C	O	0	0
			6	4	2		
2	H	1	Total	C	O	0	0
			6	4	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



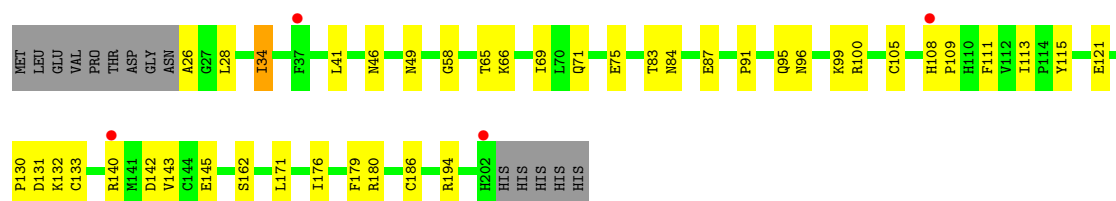
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	E	1	Total C O 4 2 2	0	0
4	H	1	Total C O 4 2 2	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	41	Total O 41 41	0	0
5	B	36	Total O 36 36	0	0
5	C	47	Total O 47 47	0	0
5	D	53	Total O 53 53	0	0
5	E	29	Total O 29 29	0	0
5	F	36	Total O 36 36	0	0
5	G	9	Total O 9 9	0	0
5	H	31	Total O 31 31	0	0

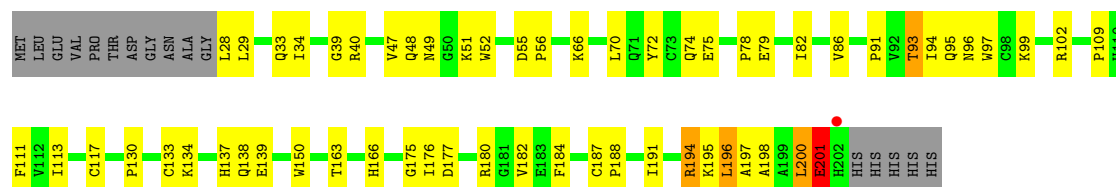
- Molecule 1: Amyloid beta A4 protein

Chain E:



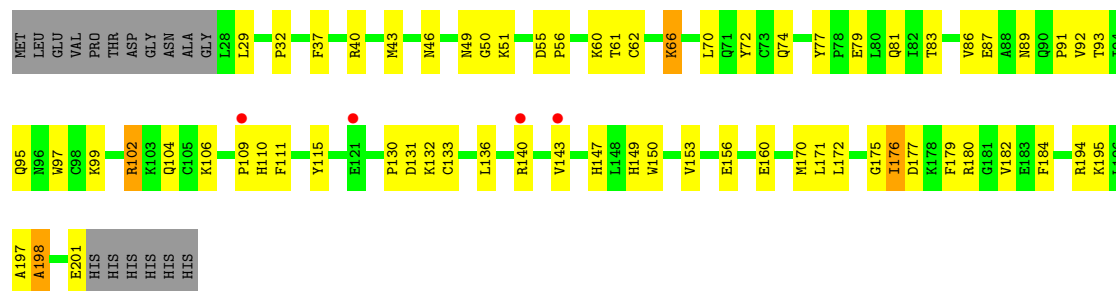
- Molecule 1: Amyloid beta A4 protein

Chain F:



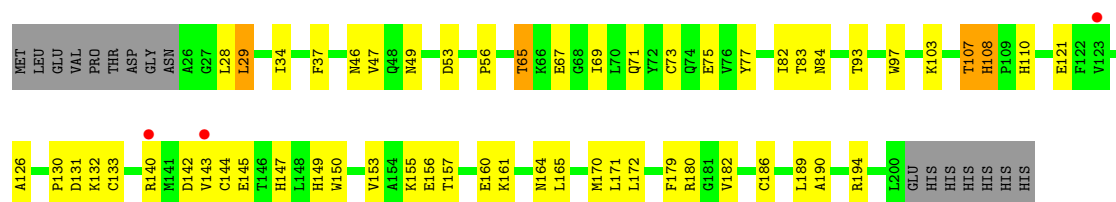
- Molecule 1: Amyloid beta A4 protein

Chain G:



- Molecule 1: Amyloid beta A4 protein

Chain H:



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	143.99Å 143.99Å 351.19Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.00 – 2.70 29.38 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.00-2.70) 99.9 (29.38-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 2.68Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.215 , 0.250 0.215 , 0.250	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	52.3	Xtriage
Anisotropy	0.079	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.7	EDS
Estimated twinning fraction	0.009 for -h,1/3*h-1/3*k-1/3*l,-4/3*h-8/3*k+1/3*l 0.009 for -1/3*h+1/3*k+1/3*l,-k,8/3*h+4/3*k+1/3*l 0.011 for -2/3*h-1/3*k-1/3*l,-1/3*h-2/3*k+1/3*l,-4/3*h+4/3*k+1/3*l 0.008 for 1/3*h+2/3*k-1/3*l,-k,-8/3*h-4/3*k-1/3*l 0.010 for -1/3*h-2/3*k+1/3*l,-2/3*h-1/3*k-1/3*l,4/3*h-4/3*k-1/3*l 0.007 for -h,2/3*h+1/3*k+1/3*l,4/3*h+8/3*k-1/3*l 0.022 for h,-h-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 74407 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11436	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.81% 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: SO4, BU4, ACT

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	0.44	0/1425	0.72	1/1929 (0.1%)
1	B	0.41	0/1401	0.68	0/1896
1	C	0.40	0/1406	0.67	0/1903
1	D	0.43	0/1405	0.70	0/1902
1	E	0.39	0/1434	0.67	0/1941
1	F	0.43	0/1425	0.70	0/1929
1	G	0.37	0/1414	0.63	0/1914
1	H	0.41	0/1414	0.67	0/1914
All	All	0.41	0/11324	0.68	1/15328 (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	A	201	GLU	N-CA-C	7.33	130.80	111.00

There are no chirality outliers.

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	1392	0	1360	41	0
1	B	1369	0	1339	50	0
1	C	1374	0	1344	54	0
1	D	1373	0	1347	25	0
1	E	1401	0	1368	27	0
1	F	1392	0	1360	51	0
1	G	1382	0	1353	57	0
1	H	1382	0	1355	56	0
2	A	6	0	10	0	0
2	B	6	0	10	1	0
2	C	6	0	10	0	0
2	D	6	0	10	1	0
2	E	6	0	10	0	0
2	F	6	0	10	2	0
2	G	6	0	10	1	0
2	H	6	0	10	1	0
3	B	5	0	0	0	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
3	E	5	0	0	0	0
3	H	5	0	0	0	0
4	B	4	0	3	1	0
4	C	4	0	3	0	0
4	E	4	0	3	0	0
4	H	4	0	3	3	0
5	A	41	0	0	1	0
5	B	36	0	0	1	0
5	C	47	0	0	3	0
5	D	53	0	0	0	0
5	E	29	0	0	0	0
5	F	36	0	0	0	0
5	G	9	0	0	0	0
5	H	31	0	0	1	0
All	All	11436	0	10918	343	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 16.

All (343) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:95:GLN:HG2	1:F:109:PRO:HG3	1.34	1.09
1:A:95:GLN:HG2	1:A:109:PRO:HG3	1.34	1.05
1:F:102:ARG:HH21	1:F:102:ARG:HB2	1.26	0.99

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:195:LYS:HE2	1:D:197:ALA:HB2	1.48	0.96
1:F:102:ARG:HB2	1:F:102:ARG:NH2	1.84	0.91
1:B:130:PRO:HG2	1:B:133:CYS:SG	2.10	0.91
1:C:146:THR:HG23	1:C:149:HIS:H	1.36	0.90
1:C:40:ARG:NH2	1:C:107:THR:HG21	1.88	0.89
1:F:195:LYS:HE2	1:F:197:ALA:HB2	1.56	0.87
1:E:95:GLN:HG2	1:E:109:PRO:HG3	1.59	0.85
1:G:61:THR:HG22	1:G:62:CYS:H	1.41	0.84
1:B:176:ILE:HD12	1:B:176:ILE:H	1.44	0.83
1:G:49:ASN:HD21	1:G:51:LYS:HB2	1.43	0.82
1:G:95:GLN:HG2	1:G:109:PRO:HG3	1.60	0.82
1:C:130:PRO:HG2	1:C:133:CYS:SG	2.19	0.81
1:A:130:PRO:HG2	1:A:133:CYS:SG	2.22	0.80
1:G:49:ASN:HD21	1:G:51:LYS:CB	1.95	0.80
1:H:73:CYS:HB3	1:H:82:ILE:HD13	1.65	0.79
1:C:140:ARG:HB2	1:C:143:VAL:HG12	1.65	0.78
1:A:103:LYS:HE2	1:C:43:MET:HE1	1.65	0.78
1:C:40:ARG:HH22	1:C:107:THR:HG21	1.48	0.75
1:H:180:ARG:HH11	1:H:180:ARG:HG2	1.51	0.75
1:C:61:THR:HG22	1:C:62:CYS:H	1.51	0.74
1:B:176:ILE:N	1:B:176:ILE:HD12	2.02	0.74
1:G:61:THR:HG22	1:G:62:CYS:N	2.04	0.73
1:G:70:LEU:O	1:G:74:GLN:HG3	1.89	0.72
1:B:91:PRO:HB3	1:B:113:ILE:HD12	1.73	0.71
1:C:143:VAL:CG2	1:C:145:GLU:HG2	2.21	0.70
1:D:95:GLN:HB3	1:D:109:PRO:HG3	1.74	0.70
1:D:130:PRO:HG2	1:D:133:CYS:SG	2.33	0.69
1:B:71:GLN:O	1:B:75:GLU:HG3	1.93	0.69
1:D:175:GLY:HA2	2:D:2:BU4:H4A	1.74	0.69
1:F:130:PRO:HG2	1:F:133:CYS:SG	2.34	0.68
1:B:102:ARG:HD2	1:D:103:LYS:O	1.93	0.68
1:C:140:ARG:HB2	1:C:143:VAL:CG1	2.24	0.67
1:E:130:PRO:HG2	1:E:133:CYS:SG	2.34	0.67
1:B:97:TRP:H	1:B:107:THR:HG22	1.60	0.67
1:A:74:GLN:HG2	1:A:82:ILE:HB	1.78	0.66
1:B:190:ALA:H	2:B:7:BU4:H2	1.61	0.65
1:F:99:LYS:HD3	1:F:102:ARG:NH2	2.12	0.65
1:H:71:GLN:O	1:H:75:GLU:HG3	1.97	0.65
1:C:143:VAL:HG22	1:C:145:GLU:HG2	1.77	0.65
1:H:143:VAL:CG2	1:H:145:GLU:HG2	2.27	0.65
1:A:103:LYS:HE2	1:C:43:MET:CE	2.28	0.64
1:A:49:ASN:OD1	1:A:51:LYS:N	2.31	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:60:LYS:HE3	1:C:75:GLU:OE2	1.96	0.64
1:G:195:LYS:HE2	1:G:197:ALA:HB2	1.79	0.64
1:G:130:PRO:HG2	1:G:133:CYS:SG	2.38	0.64
1:B:174:CYS:HB3	1:B:180:ARG:NH2	2.13	0.64
1:G:149:HIS:O	1:G:153:VAL:HG23	1.99	0.63
1:H:46:ASN:HB3	1:H:49:ASN:OD1	1.98	0.63
1:B:176:ILE:H	1:B:176:ILE:CD1	2.10	0.63
1:G:180:ARG:HH11	1:G:180:ARG:HG2	1.64	0.63
1:G:43:MET:CE	1:G:97:TRP:HA	2.29	0.63
1:F:91:PRO:HB3	1:F:113:ILE:HD12	1.81	0.63
1:B:90:GLN:OE1	4:B:208:ACT:H2	1.98	0.62
1:B:97:TRP:H	1:B:107:THR:CG2	2.13	0.62
1:G:198:ALA:HB2	1:H:126:ALA:HA	1.82	0.62
1:A:61:THR:HG22	1:A:62:CYS:N	2.13	0.62
1:A:28:LEU:HD12	1:A:47:VAL:HB	1.81	0.62
1:B:91:PRO:HB3	1:B:113:ILE:CD1	2.30	0.62
1:A:90:GLN:HB3	1:A:201:GLU:HG2	1.81	0.62
1:D:70:LEU:O	1:D:74:GLN:HG3	2.00	0.62
1:C:109:PRO:O	1:C:110:HIS:HD2	1.82	0.62
1:C:141:MET:HE3	1:F:93:THR:HB	1.80	0.62
1:G:70:LEU:HD12	1:G:70:LEU:O	2.00	0.61
1:C:61:THR:HG22	1:C:62:CYS:N	2.14	0.61
1:E:26:ALA:HB3	1:E:28:LEU:HD13	1.83	0.61
1:B:174:CYS:CB	1:B:180:ARG:NH2	2.63	0.61
1:G:46:ASN:HB3	1:G:49:ASN:OD1	2.00	0.60
1:C:164:ASN:HB2	1:C:189:LEU:HD23	1.83	0.60
1:B:174:CYS:HB3	1:B:180:ARG:HH21	1.67	0.59
1:F:150:TRP:HB3	1:F:182:VAL:HB	1.83	0.59
1:G:66:LYS:HB3	1:G:176:ILE:HG13	1.84	0.59
1:G:66:LYS:HD2	1:G:66:LYS:N	2.16	0.59
1:G:86:VAL:HA	1:G:177:ASP:OD1	2.03	0.59
1:A:61:THR:HG22	1:A:62:CYS:H	1.67	0.59
1:A:95:GLN:CG	1:A:109:PRO:HG3	2.23	0.59
1:H:164:ASN:HB2	1:H:189:LEU:HD23	1.83	0.59
1:F:137:HIS:ND1	1:F:139:GLU:OE1	2.25	0.59
1:H:46:ASN:HB2	1:H:53:ASP:OD2	2.02	0.59
1:H:133:CYS:HB3	1:H:186:CYS:O	2.03	0.59
1:H:97:TRP:H	1:H:107:THR:HG22	1.68	0.58
1:E:180:ARG:HH11	1:E:180:ARG:HG2	1.68	0.58
1:E:176:ILE:N	1:E:176:ILE:HD12	2.19	0.58
1:D:150:TRP:HB3	1:D:182:VAL:HB	1.85	0.58
1:F:70:LEU:O	1:F:74:GLN:HG3	2.04	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:71:GLN:O	1:E:75:GLU:HG3	2.04	0.57
1:E:91:PRO:HB3	1:E:113:ILE:HD13	1.86	0.57
1:G:95:GLN:CG	1:G:109:PRO:HG3	2.31	0.57
1:F:96:ASN:OD1	1:H:103:LYS:HE2	2.05	0.57
1:E:65:THR:O	1:E:69:ILE:HG13	2.05	0.57
1:G:29:LEU:HD11	1:H:71:GLN:HG2	1.87	0.57
1:H:97:TRP:H	1:H:107:THR:CG2	2.16	0.57
1:G:61:THR:CG2	1:G:62:CYS:H	2.15	0.56
1:C:146:THR:HG22	5:C:229:HOH:O	2.04	0.56
1:C:131:ASP:O	1:C:132:LYS:HB2	2.05	0.56
1:G:140:ARG:HB2	1:G:143:VAL:HG12	1.87	0.56
1:G:93:THR:HG22	1:H:180:ARG:HE	1.71	0.56
5:C:214:HOH:O	1:F:51:LYS:HG2	2.05	0.56
1:G:102:ARG:HB2	1:G:102:ARG:NH2	2.21	0.56
1:B:65:THR:O	1:B:69:ILE:HG13	2.05	0.56
1:F:86:VAL:HA	1:F:177:ASP:OD1	2.05	0.56
1:E:34:ILE:H	1:E:34:ILE:HD13	1.69	0.56
1:H:140:ARG:HB2	1:H:143:VAL:CG1	2.35	0.56
1:F:180:ARG:HG2	1:F:180:ARG:HH21	1.71	0.56
1:C:176:ILE:HD12	1:C:176:ILE:N	2.21	0.56
1:A:67:GLU:O	1:A:71:GLN:HG3	2.06	0.56
1:C:82:ILE:HD13	1:C:117:CYS:HB3	1.88	0.56
1:H:157:THR:O	1:H:161:LYS:HE2	2.05	0.55
1:H:140:ARG:HB2	1:H:143:VAL:HG12	1.87	0.55
1:G:140:ARG:CB	1:G:143:VAL:HG12	2.36	0.55
1:E:87:GLU:HG3	1:E:115:TYR:CE2	2.41	0.55
1:D:137:HIS:HD2	1:D:139:GLU:OE1	1.90	0.55
1:B:96:ASN:HA	1:B:107:THR:HG22	1.88	0.55
1:H:172:LEU:HB2	1:H:180:ARG:HB2	1.87	0.55
1:B:180:ARG:CG	1:B:180:ARG:HH21	2.20	0.55
1:C:171:LEU:O	1:C:179:PHE:HB3	2.07	0.55
1:H:144:CYS:HB2	1:H:180:ARG:HH12	1.73	0.54
1:H:67:GLU:O	1:H:71:GLN:HG3	2.07	0.54
1:A:90:GLN:CB	1:A:201:GLU:HG2	2.37	0.54
1:A:28:LEU:CD1	1:A:47:VAL:HB	2.38	0.53
1:F:49:ASN:OD1	1:F:51:LYS:N	2.40	0.53
1:D:140:ARG:HB2	1:D:143:VAL:HG13	1.89	0.53
1:B:164:ASN:HB2	1:B:189:LEU:HD23	1.90	0.53
1:E:131:ASP:O	1:E:132:LYS:HB2	2.08	0.53
1:H:155:LYS:CE	4:H:6:ACT:H3	2.39	0.53
1:E:133:CYS:HB3	1:E:186:CYS:O	2.09	0.53
1:A:70:LEU:O	1:A:74:GLN:HG3	2.09	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:155:LYS:NZ	4:H:6:ACT:H3	2.24	0.53
1:A:41:LEU:HD13	1:A:58:GLY:HA2	1.90	0.52
1:G:102:ARG:HH21	1:G:102:ARG:HB2	1.75	0.52
1:G:150:TRP:HB3	1:G:182:VAL:HB	1.92	0.52
1:H:143:VAL:HG22	1:H:145:GLU:HG2	1.91	0.52
1:C:65:THR:O	1:C:69:ILE:HG12	2.09	0.51
1:G:93:THR:CG2	1:H:180:ARG:HE	2.23	0.51
1:A:93:THR:HG23	1:A:111:PHE:CD1	2.44	0.51
1:B:67:GLU:O	1:B:71:GLN:HG3	2.11	0.51
1:G:102:ARG:HD2	1:G:104:GLN:HG3	1.93	0.51
1:G:40:ARG:HH21	1:G:106:LYS:HE2	1.76	0.51
1:G:99:LYS:HD2	1:G:102:ARG:NH1	2.25	0.51
1:A:66:LYS:HB3	1:A:176:ILE:HG22	1.92	0.51
1:C:40:ARG:CZ	1:C:99:LYS:HG3	2.41	0.51
1:A:104:GLN:NE2	1:C:103:LYS:O	2.43	0.51
1:B:131:ASP:O	1:B:132:LYS:HB2	2.11	0.50
1:C:96:ASN:HB3	1:C:105:CYS:O	2.11	0.50
1:E:66:LYS:HB3	1:E:176:ILE:HG22	1.93	0.50
1:C:133:CYS:HB3	1:C:186:CYS:O	2.12	0.50
1:G:29:LEU:HD11	1:H:71:GLN:CG	2.42	0.50
1:C:70:LEU:HD12	1:C:70:LEU:O	2.11	0.50
1:H:77:TYR:HB2	1:H:82:ILE:HD11	1.93	0.50
1:E:46:ASN:OD1	1:E:49:ASN:ND2	2.45	0.50
1:G:136:LEU:HB2	1:G:184:PHE:CZ	2.47	0.50
1:G:91:PRO:HD2	1:G:201:GLU:HG3	1.93	0.49
1:G:61:THR:CG2	1:G:62:CYS:N	2.74	0.49
1:C:73:CYS:HB3	1:C:82:ILE:HD13	1.94	0.49
1:D:191:ILE:O	1:D:194:ARG:HG2	2.13	0.49
1:F:191:ILE:CD1	1:F:196:LEU:HG	2.43	0.49
1:C:145:GLU:HG3	1:C:149:HIS:CD2	2.48	0.49
1:H:143:VAL:HG22	1:H:144:CYS:N	2.28	0.49
1:C:63:ILE:HG22	1:C:65:THR:H	1.77	0.49
1:F:33:GLN:HG3	1:F:52:TRP:CH2	2.48	0.49
1:F:200:LEU:C	1:F:201:GLU:HG2	2.33	0.49
1:B:39:GLY:O	1:B:40:ARG:HD3	2.13	0.49
1:D:163:THR:HB	1:D:187:CYS:O	2.12	0.49
1:H:29:LEU:HB2	1:H:47:VAL:O	2.12	0.49
1:B:140:ARG:HB3	1:B:143:VAL:HG23	1.95	0.49
1:A:199:ALA:C	1:A:200:LEU:HG	2.33	0.49
1:A:200:LEU:HB2	1:A:201:GLU:HG3	1.94	0.49
1:H:131:ASP:O	1:H:132:LYS:HB2	2.11	0.49
1:B:79:GLU:OE2	1:D:60:LYS:NZ	2.35	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:180:ARG:HG3	1:B:180:ARG:HH21	1.78	0.49
1:D:131:ASP:O	1:D:132:LYS:HB2	2.13	0.49
1:C:145:GLU:OE2	5:C:229:HOH:O	2.20	0.48
1:C:40:ARG:O	1:C:61:THR:HG23	2.13	0.48
1:A:195:LYS:HE2	1:A:197:ALA:HB2	1.95	0.48
1:F:175:GLY:HA2	2:F:4:BU4:H4A	1.95	0.48
1:F:66:LYS:HG2	1:F:177:ASP:HB3	1.96	0.48
1:A:93:THR:HG23	1:A:111:PHE:HD1	1.76	0.48
1:B:90:GLN:NE2	1:B:90:GLN:HA	2.27	0.48
1:C:57:SER:O	1:C:59:THR:HG23	2.13	0.48
1:C:95:GLN:HG2	1:C:109:PRO:HG3	1.94	0.48
1:C:109:PRO:O	1:C:110:HIS:CD2	2.66	0.48
1:B:172:LEU:HB2	1:B:180:ARG:HB2	1.95	0.48
1:H:34:ILE:HD12	1:H:34:ILE:C	2.34	0.48
1:H:155:LYS:HE3	4:H:6:ACT:H3	1.94	0.48
1:G:87:GLU:HG3	1:G:115:TYR:CE2	2.49	0.48
1:C:72:TYR:O	1:C:76:VAL:HG22	2.13	0.48
1:G:29:LEU:N	1:G:29:LEU:HD22	2.29	0.48
1:C:176:ILE:HD12	1:C:176:ILE:H	1.79	0.48
1:G:171:LEU:O	1:G:179:PHE:HB3	2.13	0.48
1:C:143:VAL:HG22	1:C:144:CYS:N	2.28	0.48
1:F:194:ARG:HD2	1:F:194:ARG:H	1.79	0.48
1:C:137:HIS:HD2	1:C:139:GLU:OE1	1.97	0.48
1:G:180:ARG:NH1	1:G:180:ARG:HG2	2.28	0.47
1:D:55:ASP:OD1	1:D:60:LYS:HB2	2.13	0.47
1:F:28:LEU:HG	1:F:47:VAL:HG11	1.96	0.47
1:A:30:ALA:HB3	1:A:47:VAL:HA	1.96	0.47
1:A:52:TRP:CE2	1:A:94:ILE:HD11	2.49	0.47
1:F:200:LEU:HD23	1:F:200:LEU:N	2.29	0.47
1:D:34:ILE:C	1:D:34:ILE:HD12	2.35	0.47
1:F:34:ILE:HD13	1:F:117:CYS:SG	2.55	0.47
1:A:49:ASN:OD1	1:A:50:GLY:N	2.46	0.47
1:B:150:TRP:HB3	1:B:182:VAL:HB	1.97	0.47
1:G:131:ASP:O	1:G:132:LYS:HB2	2.13	0.47
1:G:156:GLU:O	1:G:160:GLU:HG3	2.15	0.47
1:H:145:GLU:HG3	1:H:149:HIS:CD2	2.49	0.47
1:G:32:PRO:HB2	1:G:77:TYR:CE2	2.50	0.47
1:F:34:ILE:HG12	1:F:72:TYR:HD2	1.80	0.47
1:C:172:LEU:HB2	1:C:180:ARG:HB2	1.96	0.46
1:H:65:THR:O	1:H:69:ILE:HG13	2.16	0.46
1:H:130:PRO:HG2	1:H:133:CYS:SG	2.55	0.46
1:H:108:HIS:O	1:H:110:HIS:ND1	2.49	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:176:ILE:HD13	2:F:4:BU4:O3	2.16	0.46
1:H:56:PRO:HD2	5:H:12:HOH:O	2.16	0.46
1:H:140:ARG:CB	1:H:143:VAL:HG12	2.45	0.46
1:D:142:ASP:OD2	1:F:138:GLN:NE2	2.48	0.46
1:H:156:GLU:O	1:H:160:GLU:HG3	2.16	0.46
1:E:91:PRO:CB	1:E:113:ILE:HD13	2.46	0.45
1:C:42:ASN:HD21	1:C:63:ILE:CD1	2.29	0.45
1:B:143:VAL:HG12	1:B:145:GLU:HG2	1.98	0.45
1:B:96:ASN:HB3	1:B:105:CYS:O	2.16	0.45
1:A:55:ASP:HA	1:A:56:PRO:HD3	1.82	0.45
1:G:170:MET:HB3	1:G:179:PHE:CG	2.51	0.45
1:G:49:ASN:CG	1:G:50:GLY:N	2.69	0.45
1:B:83:THR:O	1:B:84:ASN:HB2	2.16	0.45
1:B:66:LYS:HB3	1:B:176:ILE:HG22	1.97	0.45
1:D:156:GLU:OE1	1:F:134:LYS:HE3	2.17	0.45
1:H:37:PHE:CE2	1:H:110:HIS:NE2	2.85	0.45
1:A:108:HIS:O	1:A:108:HIS:ND1	2.49	0.45
1:B:137:HIS:HD2	1:B:139:GLU:OE1	1.99	0.45
1:E:99:LYS:O	1:E:100:ARG:C	2.56	0.45
1:G:89:ASN:N	1:H:121:GLU:OE1	2.37	0.45
1:E:171:LEU:O	1:E:179:PHE:HB3	2.17	0.45
1:C:40:ARG:HH22	1:C:107:THR:CG2	2.25	0.45
1:E:176:ILE:CD1	1:E:176:ILE:N	2.79	0.45
1:D:93:THR:HG23	1:D:111:PHE:CD2	2.52	0.45
1:E:140:ARG:HB2	1:E:143:VAL:CG1	2.47	0.45
1:F:95:GLN:CG	1:F:109:PRO:HG3	2.25	0.44
1:H:107:THR:HG23	1:H:108:HIS:N	2.32	0.44
1:A:92:VAL:HG22	1:A:93:THR:N	2.32	0.44
1:D:144:CYS:SG	1:D:178:LYS:HB3	2.57	0.44
1:C:40:ARG:HD3	1:C:98:CYS:O	2.16	0.44
1:E:143:VAL:CG2	1:E:145:GLU:HG2	2.48	0.44
1:B:72:TYR:O	1:B:76:VAL:HG22	2.18	0.44
1:F:163:THR:HB	1:F:187:CYS:O	2.18	0.44
1:A:187:CYS:HA	1:A:188:PRO:HD3	1.83	0.44
1:G:49:ASN:ND2	1:G:51:LYS:HB2	2.23	0.44
1:H:143:VAL:CG2	1:H:144:CYS:N	2.81	0.44
1:C:149:HIS:O	1:C:153:VAL:HG23	2.18	0.44
1:H:180:ARG:CG	1:H:180:ARG:HH11	2.22	0.44
1:B:174:CYS:CB	1:B:180:ARG:HH21	2.28	0.44
1:H:34:ILE:HD12	1:H:34:ILE:O	2.18	0.44
1:B:36:MET:HG2	1:B:42:ASN:OD1	2.17	0.44
1:G:55:ASP:HA	1:G:56:PRO:HD3	1.86	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:45:MET:SD	1:B:50:GLY:O	2.75	0.44
1:G:37:PHE:CE2	1:G:110:HIS:CD2	3.05	0.44
1:H:164:ASN:ND2	1:H:165:LEU:N	2.66	0.44
1:D:140:ARG:CB	1:D:143:VAL:HG13	2.48	0.44
1:A:152:THR:O	1:A:156:GLU:HG3	2.18	0.44
1:F:191:ILE:HD12	1:F:191:ILE:N	2.33	0.44
1:G:43:MET:HE2	1:G:97:TRP:HA	2.00	0.43
1:B:97:TRP:N	1:B:107:THR:HG22	2.29	0.43
1:G:89:ASN:OD1	1:G:147:HIS:HE1	2.02	0.43
1:F:91:PRO:HB3	1:F:113:ILE:CD1	2.47	0.43
1:F:55:ASP:HA	1:F:56:PRO:HD3	1.83	0.43
1:F:99:LYS:HD3	1:F:102:ARG:HH21	1.81	0.43
1:A:61:THR:CG2	1:A:62:CYS:N	2.80	0.43
1:F:200:LEU:CD2	1:F:200:LEU:N	2.81	0.43
1:A:194:ARG:H	1:A:194:ARG:HG2	1.65	0.43
1:C:126:ALA:HA	1:F:198:ALA:HB2	2.01	0.43
1:B:77:TYR:HB3	1:B:80:LEU:HD12	2.01	0.43
1:F:187:CYS:HA	1:F:188:PRO:HD3	1.88	0.43
1:F:29:LEU:N	1:F:29:LEU:CD2	2.82	0.43
1:F:74:GLN:HG2	1:F:82:ILE:HB	2.00	0.43
1:C:63:ILE:N	1:C:63:ILE:HD12	2.34	0.42
1:A:52:TRP:CD1	1:A:94:ILE:HG12	2.54	0.42
1:H:180:ARG:NH1	1:H:180:ARG:HG2	2.26	0.42
1:B:145:GLU:HG3	1:B:149:HIS:CD2	2.54	0.42
1:F:29:LEU:HD22	1:F:29:LEU:N	2.33	0.42
1:H:149:HIS:O	1:H:153:VAL:HG23	2.19	0.42
1:H:37:PHE:CZ	1:H:110:HIS:CD2	3.07	0.42
1:A:95:GLN:HE21	1:A:109:PRO:HG3	1.84	0.42
1:H:180:ARG:CG	1:H:180:ARG:NH1	2.81	0.42
1:D:171:LEU:O	1:D:179:PHE:HB3	2.20	0.42
1:B:53:ASP:HA	5:B:218:HOH:O	2.19	0.42
1:B:90:GLN:HE21	1:B:91:PRO:HD2	1.84	0.42
1:B:109:PRO:O	1:B:110:HIS:CD2	2.72	0.42
1:B:192:GLU:OE2	1:B:192:GLU:HA	2.18	0.42
1:C:42:ASN:HD21	1:C:63:ILE:HD11	1.85	0.42
1:B:171:LEU:O	1:B:179:PHE:HB3	2.20	0.42
1:F:40:ARG:NH1	1:F:40:ARG:HG2	2.35	0.42
1:D:33:GLN:HG3	1:D:52:TRP:CH2	2.54	0.42
1:F:40:ARG:HH11	1:F:40:ARG:HG2	1.84	0.42
1:H:171:LEU:O	1:H:179:PHE:HB3	2.20	0.42
1:A:95:GLN:HE21	1:A:109:PRO:CG	2.33	0.41
1:H:190:ALA:H	2:H:8:BU4:C3	2.33	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:A:216:HOH:O	1:C:53:ASP:HA	2.20	0.41
1:C:124:SER:O	1:F:166:HIS:HE1	2.02	0.41
1:A:65:THR:O	1:A:69:ILE:HG13	2.20	0.41
1:C:143:VAL:HG21	1:C:145:GLU:HG2	1.99	0.41
1:F:33:GLN:HG3	1:F:52:TRP:HH2	1.85	0.41
1:F:39:GLY:O	1:F:40:ARG:NH1	2.53	0.41
1:H:83:THR:O	1:H:84:ASN:HB2	2.19	0.41
1:D:95:GLN:O	1:D:96:ASN:HB2	2.21	0.41
1:E:34:ILE:HD11	1:E:115:TYR:HB2	2.03	0.41
1:B:143:VAL:CG1	1:B:145:GLU:HG2	2.51	0.41
1:H:147:HIS:HD2	1:H:170:MET:CE	2.33	0.41
1:F:94:ILE:HB	1:F:97:TRP:CZ2	2.56	0.41
1:C:143:VAL:O	1:C:180:ARG:HG2	2.21	0.41
1:H:164:ASN:ND2	1:H:165:LEU:H	2.17	0.41
1:G:60:LYS:HG3	1:G:72:TYR:HE1	1.86	0.41
1:A:33:GLN:HB2	1:A:115:TYR:O	2.21	0.41
1:D:187:CYS:HA	1:D:188:PRO:HD3	1.87	0.41
1:H:179:PHE:N	1:H:179:PHE:CD1	2.89	0.41
1:G:81:GLN:O	1:G:83:THR:HG23	2.21	0.41
1:F:180:ARG:NH2	1:F:180:ARG:HG2	2.35	0.41
1:G:55:ASP:OD1	1:G:56:PRO:HD2	2.21	0.41
1:B:37:PHE:CZ	1:B:110:HIS:CE1	3.09	0.41
1:D:33:GLN:HG2	1:D:45:MET:HB2	2.02	0.41
1:H:150:TRP:HB3	1:H:182:VAL:HB	2.01	0.41
1:G:175:GLY:N	2:G:3:BU4:H2	2.35	0.41
1:C:184:PHE:N	1:C:184:PHE:CD1	2.89	0.41
1:E:83:THR:O	1:E:84:ASN:HB2	2.19	0.41
1:F:184:PHE:N	1:F:184:PHE:CD1	2.89	0.41
1:C:61:THR:CG2	1:C:62:CYS:H	2.28	0.41
1:G:92:VAL:HG22	1:G:93:THR:N	2.36	0.41
1:B:180:ARG:NH2	1:B:180:ARG:CG	2.80	0.41
1:A:61:THR:CG2	1:A:62:CYS:H	2.32	0.41
1:F:75:GLU:O	1:F:78:PRO:HD3	2.21	0.41
1:B:138:GLN:NE2	1:B:153:VAL:HG11	2.36	0.41
1:G:140:ARG:HB3	1:G:143:VAL:HG12	2.03	0.40
1:G:66:LYS:HB3	1:G:176:ILE:CG1	2.51	0.40
1:G:99:LYS:HB3	1:G:102:ARG:HB3	2.02	0.40
1:A:90:GLN:NE2	1:E:121:GLU:OE1	2.51	0.40
1:E:34:ILE:CD1	1:E:115:TYR:HB2	2.51	0.40
1:A:163:THR:HB	1:A:187:CYS:O	2.22	0.40
1:F:39:GLY:C	1:F:40:ARG:HH11	2.25	0.40
1:E:96:ASN:HB3	1:E:105:CYS:O	2.20	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:41:LEU:HD13	1:E:58:GLY:HA2	2.03	0.40
1:C:108:HIS:O	1:C:110:HIS:CD2	2.75	0.40
1:E:46:ASN:CG	1:E:49:ASN:HD21	2.25	0.40
1:B:187:CYS:HB3	1:B:188:PRO:CD	2.52	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	173/191 (91%)	164 (95%)	8 (5%)	1 (1%)	33	66
1	B	171/191 (90%)	160 (94%)	11 (6%)	0	100	100
1	C	172/191 (90%)	160 (93%)	12 (7%)	0	100	100
1	D	171/191 (90%)	165 (96%)	6 (4%)	0	100	100
1	E	175/191 (92%)	162 (93%)	13 (7%)	0	100	100
1	F	173/191 (91%)	165 (95%)	7 (4%)	1 (1%)	33	66
1	G	172/191 (90%)	161 (94%)	9 (5%)	2 (1%)	19	45
1	H	173/191 (91%)	161 (93%)	12 (7%)	0	100	100
All	All	1380/1528 (90%)	1298 (94%)	78 (6%)	4 (0%)	50	82

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	201	GLU
1	A	201	GLU
1	G	198	ALA
1	G	176	ILE

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	156/169 (92%)	150 (96%)	6 (4%)	44	76
1	B	153/169 (90%)	151 (99%)	2 (1%)	80	95
1	C	153/169 (90%)	147 (96%)	6 (4%)	43	76
1	D	154/169 (91%)	145 (94%)	9 (6%)	28	57
1	E	156/169 (92%)	150 (96%)	6 (4%)	44	76
1	F	156/169 (92%)	148 (95%)	8 (5%)	33	64
1	G	155/169 (92%)	149 (96%)	6 (4%)	43	76
1	H	154/169 (91%)	146 (95%)	8 (5%)	32	63
All	All	1237/1352 (92%)	1186 (96%)	51 (4%)	41	74

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	GLU
1	A	93	THR
1	A	111	PHE
1	A	141	MET
1	A	196	LEU
1	A	200	LEU
1	B	111	PHE
1	B	194	ARG
1	C	40	ARG
1	C	90	GLN
1	C	108	HIS
1	C	119	VAL
1	C	140	ARG
1	C	194	ARG
1	D	75	GLU
1	D	79	GLU
1	D	93	THR
1	D	109	PRO
1	D	111	PHE
1	D	139	GLU
1	D	143	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	172	LEU
1	D	194	ARG
1	E	34	ILE
1	E	108	HIS
1	E	111	PHE
1	E	142	ASP
1	E	162	SER
1	E	194	ARG
1	F	48	GLN
1	F	79	GLU
1	F	93	THR
1	F	111	PHE
1	F	194	ARG
1	F	196	LEU
1	F	200	LEU
1	F	201	GLU
1	G	66	LYS
1	G	79	GLU
1	G	102	ARG
1	G	111	PHE
1	G	172	LEU
1	G	194	ARG
1	H	28	LEU
1	H	29	LEU
1	H	65	THR
1	H	93	THR
1	H	107	THR
1	H	108	HIS
1	H	142	ASP
1	H	194	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	GLN
1	A	95	GLN
1	A	137	HIS
1	A	151	HIS
1	B	33	GLN
1	B	110	HIS
1	B	137	HIS
1	B	138	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	147	HIS
1	B	151	HIS
1	C	110	HIS
1	C	137	HIS
1	C	147	HIS
1	C	151	HIS
1	D	81	GLN
1	D	108	HIS
1	D	137	HIS
1	D	138	GLN
1	D	151	HIS
1	D	164	ASN
1	E	33	GLN
1	E	49	ASN
1	E	81	GLN
1	E	90	GLN
1	E	110	HIS
1	E	137	HIS
1	E	151	HIS
1	F	44	HIS
1	F	149	HIS
1	F	151	HIS
1	F	166	HIS
1	G	33	GLN
1	G	81	GLN
1	G	108	HIS
1	G	138	GLN
1	G	147	HIS
1	G	149	HIS
1	G	151	HIS
1	G	166	HIS
1	H	95	GLN
1	H	137	HIS
1	H	147	HIS
1	H	164	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

17 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	BU4	A	1	-	5,5,5	0.43	0	5,5,5	0.56	0
3	SO4	B	2	-	4,4,4	0.36	0	6,6,6	0.38	0
4	ACT	B	208	-	1,3,3	2.26	1 (100%)	0,3,3	0.00	-
2	BU4	B	7	-	5,5,5	0.38	0	5,5,5	0.25	0
3	SO4	C	3	-	4,4,4	0.20	0	6,6,6	0.19	0
2	BU4	C	5	-	5,5,5	0.38	0	5,5,5	0.45	0
4	ACT	C	8	-	1,3,3	2.73	1 (100%)	0,3,3	0.00	-
2	BU4	D	2	-	5,5,5	0.48	0	5,5,5	0.69	0
3	SO4	D	5	-	4,4,4	0.37	0	6,6,6	0.17	0
3	SO4	E	4	-	4,4,4	0.27	0	6,6,6	0.17	0
2	BU4	E	6	-	5,5,5	0.38	0	5,5,5	0.59	0
4	ACT	E	9	-	1,3,3	2.67	1 (100%)	0,3,3	0.00	-
2	BU4	F	4	-	5,5,5	0.52	0	5,5,5	1.28	1 (20%)
2	BU4	G	3	-	5,5,5	0.46	0	5,5,5	0.54	0
3	SO4	H	1	-	4,4,4	0.36	0	6,6,6	0.22	0
4	ACT	H	6	-	1,3,3	2.31	1 (100%)	0,3,3	0.00	-
2	BU4	H	8	-	5,5,5	0.40	0	5,5,5	0.94	1 (20%)

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	BU4	A	1	-	-	0/3/3/3	0/0/0/0
3	SO4	B	2	-	-	0/0/0/0	0/0/0/0
4	ACT	B	208	-	-	0/0/0/0	0/0/0/0
2	BU4	B	7	-	-	0/3/3/3	0/0/0/0
3	SO4	C	3	-	-	0/0/0/0	0/0/0/0
2	BU4	C	5	-	-	0/3/3/3	0/0/0/0
4	ACT	C	8	-	-	0/0/0/0	0/0/0/0
2	BU4	D	2	-	-	0/3/3/3	0/0/0/0
3	SO4	D	5	-	-	0/0/0/0	0/0/0/0
3	SO4	E	4	-	-	0/0/0/0	0/0/0/0
2	BU4	E	6	-	-	0/3/3/3	0/0/0/0
4	ACT	E	9	-	-	0/0/0/0	0/0/0/0
2	BU4	F	4	-	-	0/3/3/3	0/0/0/0
2	BU4	G	3	-	-	0/3/3/3	0/0/0/0
3	SO4	H	1	-	-	0/0/0/0	0/0/0/0
4	ACT	H	6	-	-	0/0/0/0	0/0/0/0
2	BU4	H	8	-	-	0/3/3/3	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	8	ACT	CH3-C	2.73	1.52	1.48
4	E	9	ACT	CH3-C	2.67	1.52	1.48
4	H	6	ACT	CH3-C	2.31	1.52	1.48
4	B	208	ACT	CH3-C	2.26	1.52	1.48

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	4	BU4	C1-C2-C3	2.82	120.67	113.40
2	H	8	BU4	C1-C2-C3	2.03	118.63	113.40

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	175/191 (91%)	-0.29	2 (1%) 77 82	19, 41, 72, 86	14 (8%)
1	B	173/191 (90%)	-0.23	1 (0%) 86 90	26, 45, 66, 89	10 (5%)
1	C	174/191 (91%)	-0.23	1 (0%) 86 90	32, 46, 72, 87	10 (5%)
1	D	173/191 (90%)	-0.33	2 (1%) 75 81	21, 38, 66, 81	14 (8%)
1	E	177/191 (92%)	-0.18	4 (2%) 57 64	31, 49, 79, 100	16 (9%)
1	F	175/191 (91%)	-0.27	1 (0%) 86 90	25, 44, 75, 97	15 (8%)
1	G	174/191 (91%)	0.06	4 (2%) 57 64	36, 60, 89, 100	22 (12%)
1	H	175/191 (91%)	-0.23	3 (1%) 67 73	29, 48, 72, 89	17 (9%)
All	All	1396/1528 (91%)	-0.21	18 (1%) 74 79	19, 47, 75, 100	118 (8%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	202	HIS	5.4
1	D	29	LEU	3.5
1	G	143	VAL	3.5
1	F	202	HIS	3.3
1	B	108	HIS	3.2
1	H	143	VAL	2.9
1	E	108	HIS	2.7
1	D	109	PRO	2.5
1	G	140	ARG	2.5
1	E	37	PHE	2.5
1	H	123	VAL	2.5
1	G	121	GLU	2.5
1	H	140	ARG	2.4
1	A	202	HIS	2.3
1	E	140	ARG	2.3
1	C	105	CYS	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	G	109	PRO	2.1
1	A	109	PRO	2.1

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
4	ACT	C	8	4/4	0.42	45.14	85,86,87,88	0
4	ACT	E	9	4/4	0.27	12.39	71,72,73,74	0
4	ACT	B	208	4/4	0.37	10.90	68,68,70,73	0
4	ACT	H	6	4/4	0.33	8.95	57,57,58,60	0
2	BU4	F	4	6/6	0.26	5.60	62,70,76,76	0
2	BU4	G	3	6/6	0.28	4.21	78,82,83,85	0
2	BU4	D	2	6/6	0.18	2.08	59,60,66,66	0
2	BU4	A	1	6/6	0.17	1.66	50,53,55,55	0
3	SO4	C	3	5/5	0.17	1.28	38,42,49,49	0
3	SO4	E	4	5/5	0.17	1.21	39,45,48,53	0
2	BU4	H	8	6/6	0.22	0.97	50,57,65,67	0
2	BU4	E	6	6/6	0.20	0.63	63,65,71,75	0
2	BU4	B	7	6/6	0.16	0.34	58,64,65,65	0
3	SO4	B	2	5/5	0.15	-0.45	27,36,46,50	0
2	BU4	C	5	6/6	0.11	-0.75	51,53,58,58	0
3	SO4	H	1	5/5	0.12	-0.79	39,42,48,50	0
3	SO4	D	5	5/5	0.09	-1.07	72,73,75,75	0

6.5 Other polymers ⓘ

There are no such residues in this entry.