



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 12:33 AM GMT

PDB ID : 3I3B
Title : E.coli (lacZ) Beta-Galactosidase (M542A) in Complex with D-Galactopyranosyl-1-on
Authors : Dugdale, M.L.; Dymianiw, D.; Minhas, B.; Huber, R.E.
Deposited on : 2009-06-30
Resolution : 2.20 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>

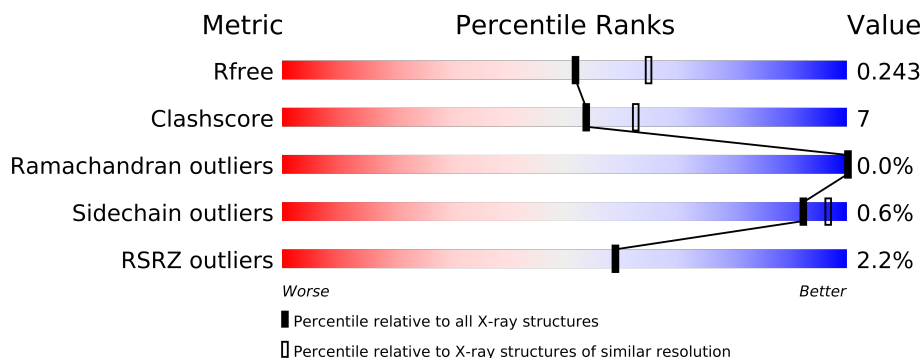
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.20 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}	66092	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (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. 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	1023	
1	B	1023	
1	C	1023	
1	D	1023	

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	149	A	2001	-	X
2	149	D	2001	-	X
4	NA	D	3104	-	X
5	DMS	A	5002	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
5	DMS	A	5004	-	X
5	DMS	A	5006	-	X
5	DMS	A	5007	-	X
5	DMS	A	5009	-	X
5	DMS	A	5010	-	X
5	DMS	A	5014	-	X
5	DMS	A	5015	-	X
5	DMS	A	5016	-	X
5	DMS	A	5017	-	X
5	DMS	A	5019	-	X
5	DMS	A	5024	-	X
5	DMS	A	5125	-	X
5	DMS	A	5127	-	X
5	DMS	B	5002	-	X
5	DMS	B	5003	-	X
5	DMS	B	5004	-	X
5	DMS	B	5006	-	X
5	DMS	B	5007	-	X
5	DMS	B	5009	-	X
5	DMS	B	5011	-	X
5	DMS	B	5013	-	X
5	DMS	B	5015	-	X
5	DMS	B	5017	-	X
5	DMS	B	5018	-	X
5	DMS	B	5019	-	X
5	DMS	B	5020	-	X
5	DMS	B	5021	-	X
5	DMS	B	5022	-	X
5	DMS	B	5025	-	X
5	DMS	B	5026	-	X
5	DMS	B	5027	-	X
5	DMS	B	5028	-	X
5	DMS	B	5131	-	X
5	DMS	B	5132	-	X
5	DMS	C	1024	-	X
5	DMS	C	5003	-	X
5	DMS	C	5004	-	X
5	DMS	C	5006	-	X
5	DMS	C	5008	-	X
5	DMS	C	5009	-	X
5	DMS	C	5012	-	X
5	DMS	C	5013	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
5	DMS	C	5014	-	X
5	DMS	C	5016	-	X
5	DMS	C	5017	-	X
5	DMS	C	5018	-	X
5	DMS	C	5019	-	X
5	DMS	C	5020	-	X
5	DMS	C	5021	-	X
5	DMS	C	5022	-	X
5	DMS	C	5027	-	X
5	DMS	C	5028	-	X
5	DMS	C	5032	-	X
5	DMS	C	5033	-	X
5	DMS	C	5134	-	X
5	DMS	D	5002	-	X
5	DMS	D	5004	-	X
5	DMS	D	5007	-	X
5	DMS	D	5010	-	X
5	DMS	D	5013	-	X
5	DMS	D	5014	-	X
5	DMS	D	5015	-	X
5	DMS	D	5018	-	X
5	DMS	D	5019	-	X
5	DMS	D	5020	-	X
5	DMS	D	5021	-	X
5	DMS	D	5022	-	X
5	DMS	D	5023	-	X
5	DMS	D	5024	-	X
5	DMS	D	5026	-	X
5	DMS	D	5027	-	X
5	DMS	D	5229	-	X
5	DMS	D	5230	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 36275 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 Beta-galactosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1011	Total	C	N	O	S	0	0	0
			8122	5136	1440	1509	37			
1	B	1011	Total	C	N	O	S	0	0	0
			8122	5136	1440	1509	37			
1	C	1011	Total	C	N	O	S	0	0	0
			8122	5136	1440	1509	37			
1	D	1011	Total	C	N	O	S	0	0	0
			8122	5136	1440	1509	37			

There are 36 discrepancies between the modelled and reference sequences:

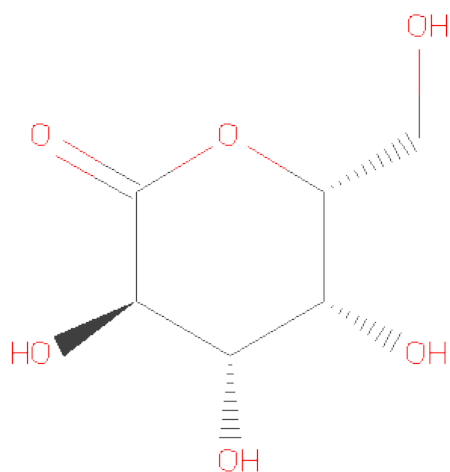
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	EXPRESSION TAG	UNP B8LFD6
A	2	SER	-	EXPRESSION TAG	UNP B8LFD6
A	3	HIS	-	EXPRESSION TAG	UNP B8LFD6
A	4	MET	-	EXPRESSION TAG	UNP B8LFD6
A	5	LEU	-	EXPRESSION TAG	UNP B8LFD6
A	6	GLU	-	EXPRESSION TAG	UNP B8LFD6
A	7	ASP	-	EXPRESSION TAG	UNP B8LFD6
A	8	PRO	-	EXPRESSION TAG	UNP B8LFD6
A	542	ALA	MET	ENGINEERED	UNP B8LFD6
B	1	GLY	-	EXPRESSION TAG	UNP B8LFD6
B	2	SER	-	EXPRESSION TAG	UNP B8LFD6
B	3	HIS	-	EXPRESSION TAG	UNP B8LFD6
B	4	MET	-	EXPRESSION TAG	UNP B8LFD6
B	5	LEU	-	EXPRESSION TAG	UNP B8LFD6
B	6	GLU	-	EXPRESSION TAG	UNP B8LFD6
B	7	ASP	-	EXPRESSION TAG	UNP B8LFD6
B	8	PRO	-	EXPRESSION TAG	UNP B8LFD6
B	542	ALA	MET	ENGINEERED	UNP B8LFD6
C	1	GLY	-	EXPRESSION TAG	UNP B8LFD6
C	2	SER	-	EXPRESSION TAG	UNP B8LFD6
C	3	HIS	-	EXPRESSION TAG	UNP B8LFD6

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Chain	Residue	Modelled	Actual	Comment	Reference
C	4	MET	-	EXPRESSION TAG	UNP B8LFD6
C	5	LEU	-	EXPRESSION TAG	UNP B8LFD6
C	6	GLU	-	EXPRESSION TAG	UNP B8LFD6
C	7	ASP	-	EXPRESSION TAG	UNP B8LFD6
C	8	PRO	-	EXPRESSION TAG	UNP B8LFD6
C	542	ALA	MET	ENGINEERED	UNP B8LFD6
D	1	GLY	-	EXPRESSION TAG	UNP B8LFD6
D	2	SER	-	EXPRESSION TAG	UNP B8LFD6
D	3	HIS	-	EXPRESSION TAG	UNP B8LFD6
D	4	MET	-	EXPRESSION TAG	UNP B8LFD6
D	5	LEU	-	EXPRESSION TAG	UNP B8LFD6
D	6	GLU	-	EXPRESSION TAG	UNP B8LFD6
D	7	ASP	-	EXPRESSION TAG	UNP B8LFD6
D	8	PRO	-	EXPRESSION TAG	UNP B8LFD6
D	542	ALA	MET	ENGINEERED	UNP B8LFD6

- Molecule 2 is SUGAR (D-GALACTONOLACTONE) (three-letter code: 149) (formula: $C_6H_{10}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			12	6	6		
2	B	1	Total	C	O	0	0
			12	6	6		
2	C	1	Total	C	O	0	0
			12	6	6		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			12	6	6		

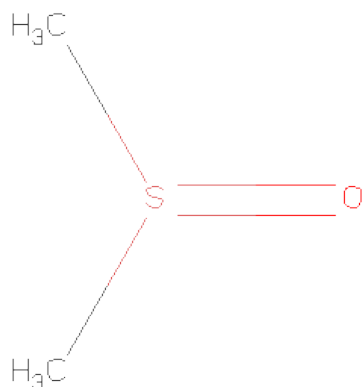
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	3	Total	Mg	0	0
			3	3		
3	A	4	Total	Mg	0	0
			4	4		
3	D	2	Total	Mg	0	0
			2	2		
3	C	2	Total	Mg	0	0
			2	2		

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	4	Total	Na	0	0
			4	4		
4	A	4	Total	Na	0	0
			4	4		
4	D	4	Total	Na	0	0
			4	4		
4	C	4	Total	Na	0	0
			4	4		

- Molecule 5 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C₂H₆OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	O	S	0	0
			4	2	1	1		
5	A	1	Total	C	O	S	0	0
			4	2	1	1		
5	A	1	Total	C	O	S	0	0
			4	2	1	1		
5	A	1	Total	C	O	S	0	0
			4	2	1	1		
5	A	1	Total	C	O	S	0	0
			4	2	1	1		
5	A	1	Total	C	O	S	0	0
			4	2	1	1		
5	A	1	Total	C	O	S	0	0
			4	2	1	1		
5	A	1	Total	C	O	S	0	0
			4	2	1	1		
5	A	1	Total	C	O	S	0	0
			4	2	1	1		
5	A	1	Total	C	O	S	0	0
			4	2	1	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	A	1	Total 4	C 2	O 1	S 1	0	0
5	B	1	Total 4	C 2	O 1	S 1	0	0
5	B	1	Total 4	C 2	O 1	S 1	0	0
5	B	1	Total 4	C 2	O 1	S 1	0	0
5	B	1	Total 4	C 2	O 1	S 1	0	0
5	B	1	Total 4	C 2	O 1	S 1	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total 4	C 2	O 1	S 1	0	0
5	B	1	Total 4	C 2	O 1	S 1	0	0
5	B	1	Total 4	C 2	O 1	S 1	0	0
5	B	1	Total 4	C 2	O 1	S 1	0	0
5	B	1	Total 4	C 2	O 1	S 1	0	0
5	B	1	Total 4	C 2	O 1	S 1	0	0
5	B	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	C	1	Total 4	C 2	O 1	S 1	0	0
5	D	1	Total 4	C 2	O 1	S 1	0	0
5	D	1	Total 4	C 2	O 1	S 1	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	D	1	Total 4	C 2	O 1	S 1	0	0
5	D	1	Total 4	C 2	O 1	S 1	0	0
5	D	1	Total 4	C 2	O 1	S 1	0	0
5	D	1	Total 4	C 2	O 1	S 1	0	0
5	D	1	Total 4	C 2	O 1	S 1	0	0
5	D	1	Total 4	C 2	O 1	S 1	0	0
5	D	1	Total 4	C 2	O 1	S 1	0	0

- Molecule 6 is water.

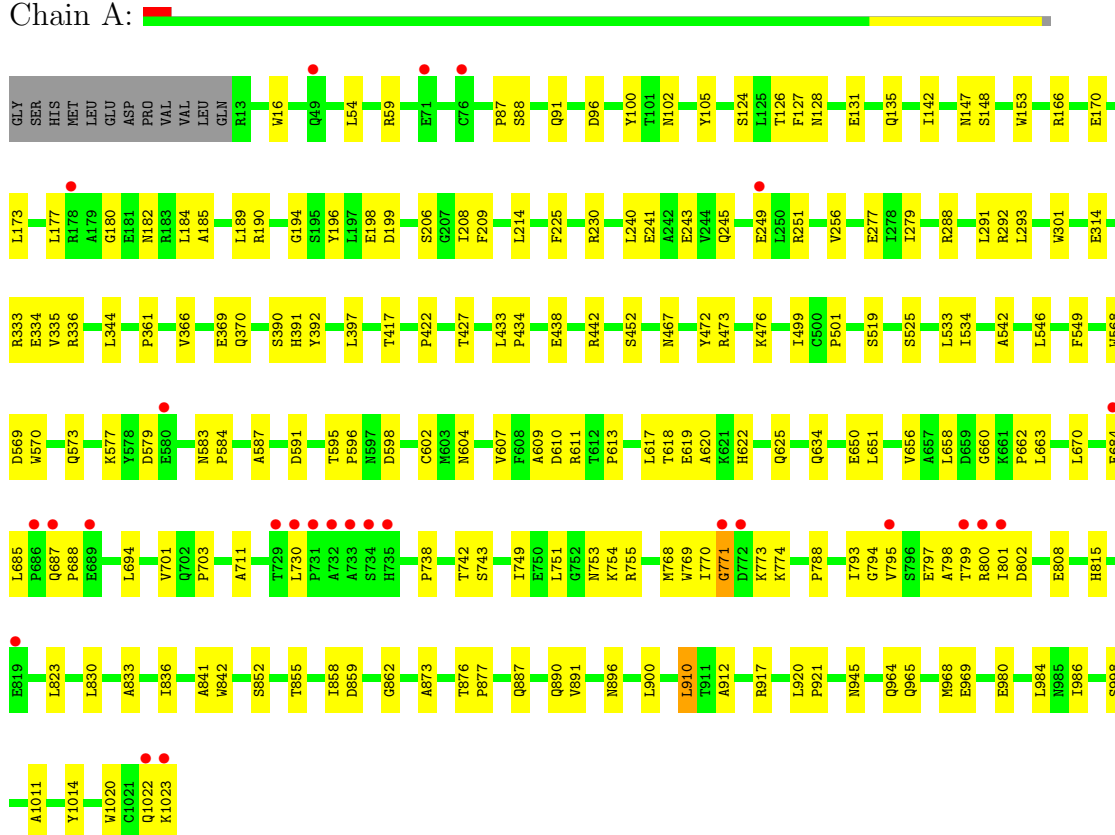
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	798	Total 798	O 798	0	0
6	B	821	Total 821	O 821	0	0
6	C	785	Total 785	O 785	0	0
6	D	804	Total 804	O 804	0	0

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.

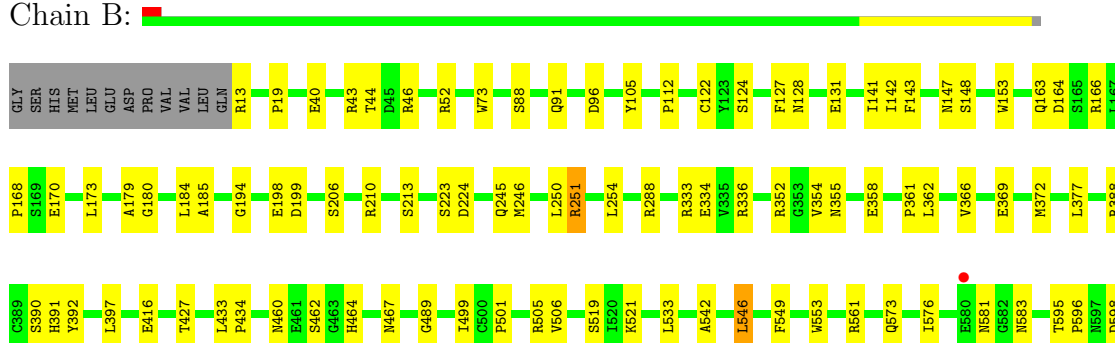
• Molecule 1: Beta-galactosidase

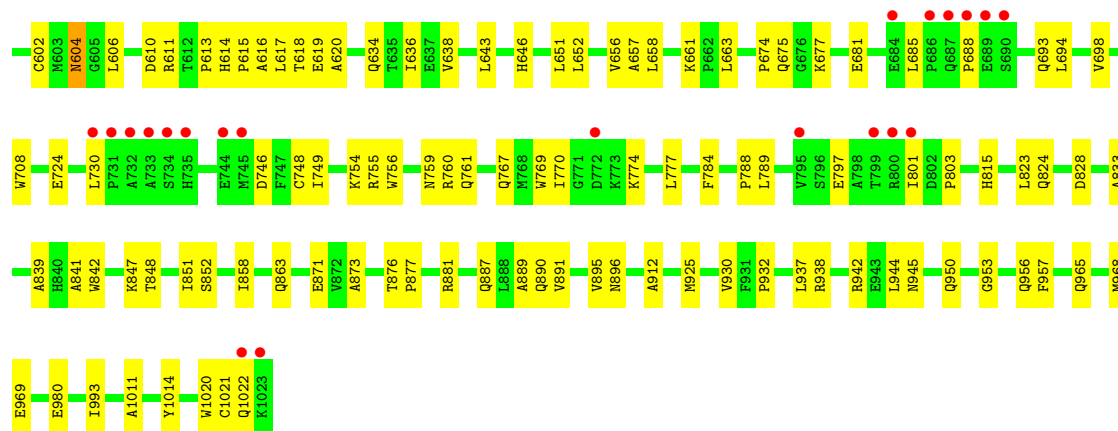
Chain A:



• Molecule 1: Beta-galactosidase

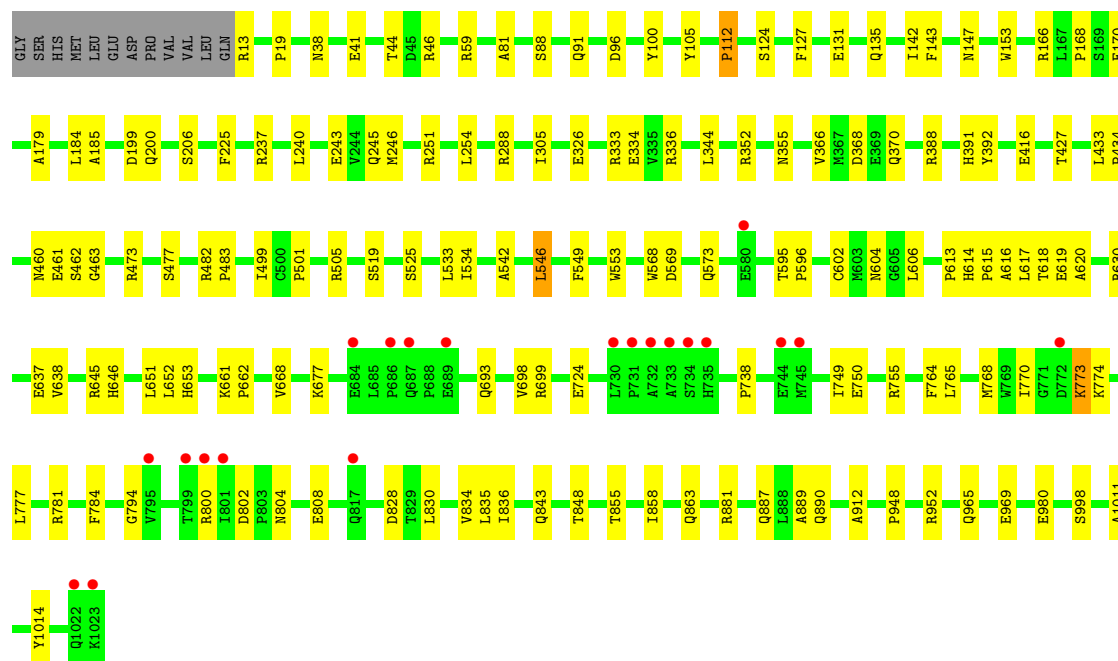
Chain B:





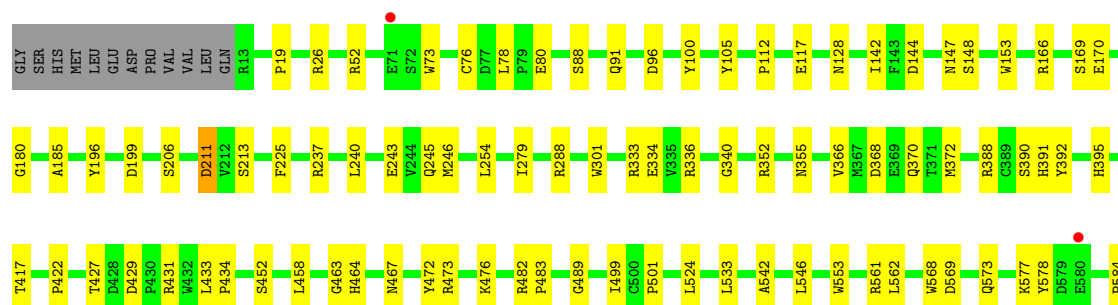
• Molecule 1: Beta-galactosidase

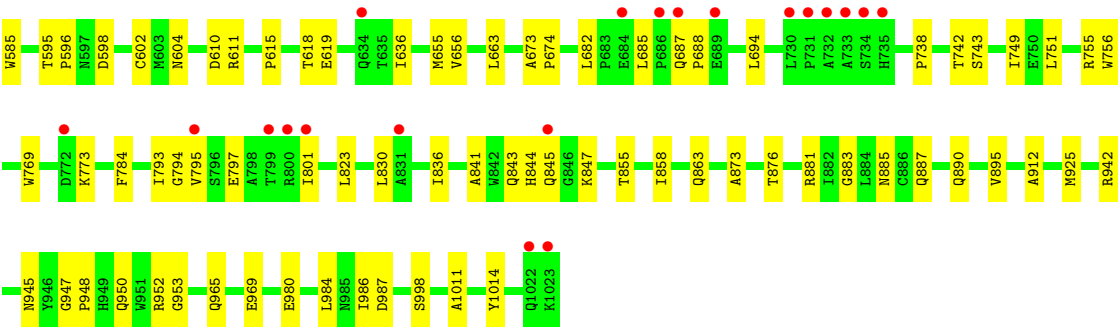
Chain C:



• Molecule 1: Beta-galactosidase

Chain D:





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	149.60Å 168.42Å 201.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.18 – 2.20 14.18 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.7 (14.18-2.20) 96.2 (14.18-2.00)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	0.17	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 2.00Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.193 , 0.244 0.191 , 0.243	Depositor DCC
R_{free} test set	3517 reflections (1.44%)	DCC
Wilson B-factor (Å ²)	19.6	Xtriage
Anisotropy	0.381	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 46.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 326526 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	36275	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 36.26 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 5.1383e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: NA, MG, DMS, 149

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	0.33	0/8364	0.62	1/11412 (0.0%)
1	B	0.33	0/8364	0.63	0/11412
1	C	0.33	0/8364	0.62	1/11412 (0.0%)
1	D	0.33	0/8364	0.62	0/11412
All	All	0.33	0/33456	0.62	2/45648 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	771	GLY	N-CA-C	-5.33	99.76	113.10
1	C	770	ILE	N-CA-C	-5.08	97.28	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	8122	0	7711	130	0
1	B	8122	0	7712	125	0
1	C	8122	0	7712	102	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	8122	0	7712	104	0
2	A	12	0	9	0	0
2	B	12	0	9	0	0
2	C	12	0	9	0	0
2	D	12	0	9	0	0
3	A	4	0	0	0	0
3	B	3	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
4	A	4	0	0	0	0
4	B	4	0	0	0	0
4	C	4	0	0	0	0
4	D	4	0	0	0	0
5	A	120	0	180	2	0
5	B	132	0	198	0	0
5	C	132	0	198	1	0
5	D	120	0	180	0	0
6	A	798	0	0	2	0
6	B	821	0	0	4	0
6	C	785	0	0	2	0
6	D	804	0	0	4	0
All	All	36275	0	31639	449	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 7.

All (449) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:142:ILE:HG12	1:A:170:GLU:HG2	1.42	1.01
1:A:249:GLU:CD	1:A:251:ARG:HE	1.68	0.96
1:C:142:ILE:HG12	1:C:170:GLU:HG2	1.59	0.82
1:A:730:LEU:HD21	1:B:823:LEU:O	1.82	0.80
1:A:147:ASN:HB3	1:A:206:SER:HA	1.63	0.79
1:B:142:ILE:HG12	1:B:170:GLU:HG2	1.63	0.79
1:B:767:GLN:NE2	1:B:774:LYS:HB3	2.00	0.77
1:B:245:GLN:HG2	1:B:288:ARG:HG2	1.69	0.75
1:C:773:LYS:HE2	1:C:774:LYS:N	2.01	0.75
1:D:237:ARG:HH11	1:D:237:ARG:HB3	1.53	0.74
1:D:153:TRP:HB2	1:D:185:ALA:HB3	1.69	0.74
1:D:237:ARG:NH1	1:D:237:ARG:HB3	2.03	0.74
1:B:863:GLN:HG2	1:B:1021:CYS:HB3	1.69	0.74
1:C:44:THR:OG1	1:C:46:ARG:HD3	1.88	0.73

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:88:SER:HA	1:D:366:VAL:HG21	1.71	0.72
1:D:863:GLN:HE22	1:D:952:ARG:HH21	1.37	0.70
1:A:241:GLU:HG2	1:A:292:ARG:HG2	1.73	0.70
1:A:245:GLN:HG2	1:A:288:ARG:HG2	1.74	0.70
1:D:147:ASN:HB3	1:D:206:SER:HA	1.75	0.68
1:B:52:ARG:O	1:B:213:SER:HB2	1.94	0.68
1:C:245:GLN:HG2	1:C:288:ARG:HG2	1.77	0.67
1:A:945:ASN:HB3	1:A:1023:LYS:NZ	2.11	0.66
1:B:890:GLN:HG2	1:B:891:VAL:N	2.12	0.65
1:D:965:GLN:O	1:D:969:GLU:HG3	1.95	0.65
1:A:54:LEU:HD11	1:A:214:LEU:HD13	1.78	0.65
1:C:473:ARG:HH12	1:C:477:SER:HB2	1.62	0.65
1:C:1011:ALA:HB3	1:C:1014:TYR:CZ	2.33	0.64
1:A:579:ASP:OD2	1:A:583:ASN:HB2	1.97	0.64
1:C:237:ARG:NH1	1:C:237:ARG:HB3	2.12	0.64
1:D:749:ILE:N	1:D:749:ILE:HD12	2.13	0.64
1:A:88:SER:HA	1:A:366:VAL:HG21	1.79	0.64
1:A:788:PRO:HD2	1:A:968:MET:HG3	1.81	0.63
1:C:147:ASN:HB3	1:C:206:SER:HA	1.79	0.63
1:A:749:ILE:HD12	1:A:858:ILE:HD12	1.81	0.63
1:A:249:GLU:HG2	1:A:251:ARG:HG2	1.79	0.62
1:D:372:MET:HE1	1:D:395:HIS:HB3	1.80	0.62
1:D:844:HIS:ND1	1:D:845:GLN:HG2	2.14	0.62
1:C:305:ILE:HD11	1:C:645:ARG:HB3	1.79	0.62
1:C:131:GLU:O	1:C:135:GLN:HG3	1.99	0.62
1:A:823:LEU:O	1:B:730:LEU:HD11	2.00	0.62
1:D:578:TYR:CE1	1:D:584:PRO:HB3	2.35	0.62
1:C:91:GLN:HG3	1:C:96:ASP:OD1	2.00	0.61
1:D:573:GLN:HB2	1:D:602:CYS:O	2.01	0.61
1:B:777:LEU:HG	1:B:889:ALA:HA	1.82	0.61
1:C:355:ASN:OD1	1:C:388:ARG:HD3	2.01	0.60
1:D:863:GLN:HE22	1:D:952:ARG:NH2	1.99	0.60
1:A:801:ILE:N	1:A:801:ILE:HD12	2.16	0.60
1:B:334:GLU:OE1	1:B:336:ARG:NH1	2.33	0.60
1:A:751:LEU:HD23	1:A:862:GLY:HA2	1.83	0.60
1:B:873:ALA:O	1:B:876:THR:HG22	2.01	0.59
1:C:887:GLN:NE2	1:C:980:GLU:O	2.35	0.59
1:B:1011:ALA:HB3	1:B:1014:TYR:CZ	2.37	0.59
1:A:965:GLN:O	1:A:969:GLU:HG3	2.02	0.59
1:C:473:ARG:NH1	1:C:477:SER:HB2	2.18	0.59
1:B:355:ASN:OD1	1:B:388:ARG:HD3	2.01	0.59
1:B:643:LEU:HD23	1:B:675:GLN:NE2	2.18	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:254:LEU:HD12	6:D:4828:HOH:O	2.03	0.59
1:C:131:GLU:OE1	1:C:179:ALA:HB2	2.03	0.59
1:B:774:LYS:HB2	6:B:4851:HOH:O	2.03	0.59
1:D:433:LEU:HB3	1:D:434:PRO:HD3	1.85	0.58
1:D:793:ILE:HG22	1:D:795:VAL:HG22	1.84	0.58
1:C:166:ARG:HG3	1:C:392:TYR:HB2	1.84	0.58
1:A:131:GLU:O	1:A:135:GLN:HG2	2.03	0.58
1:A:153:TRP:HB2	1:A:185:ALA:HB3	1.83	0.58
1:D:656:VAL:HG21	1:D:685:LEU:HD13	1.84	0.58
1:C:777:LEU:HG	1:C:889:ALA:HA	1.86	0.58
1:B:754:LYS:HE2	1:B:1022:GLN:OE1	2.04	0.58
1:B:754:LYS:HE3	1:B:770:ILE:HG23	1.85	0.58
1:B:377:LEU:CD2	1:B:708:TRP:HA	2.34	0.58
1:A:334:GLU:OE1	1:A:336:ARG:NH1	2.37	0.58
1:C:749:ILE:HD12	1:C:749:ILE:N	2.18	0.58
1:A:249:GLU:CG	1:A:251:ARG:HE	2.17	0.57
1:D:245:GLN:HG2	1:D:288:ARG:HG2	1.85	0.57
1:A:613:PRO:HB3	1:A:617:LEU:HD23	1.86	0.57
1:B:890:GLN:HG2	1:B:891:VAL:H	1.69	0.57
1:B:651:LEU:HD12	1:B:651:LEU:C	2.25	0.57
1:C:533:LEU:C	1:C:533:LEU:HD23	2.25	0.56
1:C:200:GLN:HG2	1:C:391:HIS:HB2	1.87	0.56
1:A:279:ILE:HD11	1:D:422:PRO:HG3	1.86	0.56
1:B:823:LEU:HD11	1:B:841:ALA:HB2	1.88	0.56
1:A:194:GLY:O	1:A:198:GLU:HG3	2.06	0.56
1:D:615:PRO:O	1:D:618:THR:HG22	2.06	0.56
1:B:759:ASN:OD1	1:B:761:GLN:HB3	2.06	0.56
1:A:823:LEU:HD11	1:A:841:ALA:HB2	1.86	0.56
1:B:131:GLU:OE1	1:B:179:ALA:HB2	2.05	0.56
1:B:499:ILE:HG22	1:B:501:PRO:HD3	1.88	0.55
1:B:746:ASP:HA	1:B:760:ARG:HG3	1.88	0.55
1:B:769:TRP:NE1	1:B:774:LYS:HG2	2.22	0.55
1:D:863:GLN:NE2	1:D:952:ARG:HH21	2.04	0.55
1:C:81:ALA:HA	5:C:5030:DMS:H11	1.89	0.55
1:A:96:ASP:OD2	1:A:190:ARG:NH1	2.40	0.55
1:B:965:GLN:O	1:B:969:GLU:HG3	2.06	0.55
1:B:749:ILE:N	1:B:749:ILE:HD12	2.21	0.55
1:B:573:GLN:HB2	1:B:602:CYS:O	2.06	0.55
1:A:634:GLN:NE2	1:A:684:GLU:HA	2.21	0.55
1:D:499:ILE:HG22	1:D:501:PRO:HD3	1.88	0.55
1:B:688:PRO:HD3	1:B:694:LEU:HD11	1.89	0.55
1:D:237:ARG:HH11	1:D:237:ARG:CB	2.17	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:88:SER:HA	1:C:366:VAL:HG21	1.90	0.54
1:A:472:TYR:O	1:A:476:LYS:HG2	2.07	0.54
1:C:863:GLN:NE2	1:C:952:ARG:NH2	2.55	0.54
1:B:88:SER:HA	1:B:366:VAL:HG21	1.87	0.54
1:B:784:PHE:HA	1:B:881:ARG:O	2.08	0.54
1:C:334:GLU:OE1	1:C:336:ARG:NH1	2.39	0.54
1:B:851:ILE:HB	1:B:871:GLU:HB2	1.90	0.53
1:D:890:GLN:HE22	1:D:948:PRO:HD3	1.73	0.53
1:D:945:ASN:OD1	1:D:950:GLN:HG3	2.09	0.53
1:A:105:TYR:CE1	1:A:199:ASP:HB2	2.44	0.53
1:A:127:PHE:CE1	1:A:184:LEU:HG	2.43	0.53
1:C:750:GLU:OE2	1:C:755:ARG:HD2	2.09	0.53
1:C:773:LYS:HE2	1:C:774:LYS:H	1.72	0.53
1:D:823:LEU:HD11	1:D:841:ALA:HB2	1.91	0.53
1:B:652:LEU:HD11	1:B:698:VAL:HB	1.91	0.53
1:B:127:PHE:HE1	1:B:184:LEU:HG	1.74	0.53
1:A:984:LEU:HD21	1:A:986:ILE:HD11	1.91	0.53
1:A:663:LEU:HD11	1:A:688:PRO:HB3	1.91	0.53
1:C:352:ARG:HG2	1:C:553:TRP:CH2	2.43	0.53
1:B:887:GLN:NE2	1:B:980:GLU:O	2.40	0.53
1:C:863:GLN:HE22	1:C:952:ARG:NH2	2.07	0.53
1:D:73:TRP:HB2	1:D:78:LEU:HD11	1.91	0.53
1:A:196:TYR:O	1:A:417:THR:HG22	2.09	0.53
1:B:606:LEU:O	1:B:614:HIS:HB2	2.09	0.52
1:D:887:GLN:NE2	1:D:980:GLU:O	2.42	0.52
1:B:369:GLU:HA	1:B:372:MET:HE3	1.91	0.52
1:C:646:HIS:HB3	6:C:4937:HOH:O	2.09	0.52
1:B:755:ARG:HD3	1:B:769:TRP:CE3	2.45	0.52
1:C:843:GLN:HG2	1:C:848:THR:HA	1.91	0.52
1:A:549:PHE:CE2	1:A:620:ALA:HA	2.45	0.52
1:C:499:ILE:HG22	1:C:501:PRO:HD3	1.92	0.52
1:D:473:ARG:O	1:D:473:ARG:NH1	2.42	0.52
1:C:127:PHE:CE1	1:C:184:LEU:HG	2.45	0.52
1:B:542:ALA:HA	1:B:604:ASN:HA	1.91	0.52
1:B:433:LEU:HB3	1:B:434:PRO:HD3	1.92	0.52
1:B:595:THR:HA	1:B:596:PRO:C	2.29	0.52
1:A:873:ALA:O	1:A:876:THR:HG22	2.09	0.52
1:C:237:ARG:CB	1:C:237:ARG:HH11	2.22	0.51
1:B:777:LEU:HD11	1:B:980:GLU:HG2	1.92	0.51
1:B:863:GLN:HG2	1:B:1021:CYS:CB	2.38	0.51
1:A:945:ASN:HB3	1:A:1023:LYS:HZ2	1.75	0.51
1:D:19:PRO:HD3	1:D:112:PRO:CB	2.41	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:251:ARG:HB2	1:C:254:LEU:HG	1.92	0.51
1:B:369:GLU:HG3	1:B:397:LEU:HD21	1.93	0.51
1:C:240:LEU:HD23	1:C:240:LEU:C	2.31	0.51
1:B:73:TRP:CE2	1:B:122:CYS:HB3	2.45	0.51
1:C:237:ARG:HH11	1:C:237:ARG:HB3	1.75	0.51
1:D:105:TYR:CE1	1:D:199:ASP:HB2	2.45	0.51
1:D:883:GLY:HA3	1:D:987:ASP:HA	1.93	0.51
1:B:427:THR:HG21	1:B:462:SER:HB3	1.93	0.51
1:B:937:LEU:HA	1:B:957:PHE:O	2.11	0.51
1:B:930:VAL:O	1:B:932:PRO:HD3	2.11	0.50
1:A:749:ILE:CD1	1:A:858:ILE:HD12	2.41	0.50
1:D:595:THR:HA	1:D:596:PRO:C	2.32	0.50
1:D:91:GLN:HG3	1:D:96:ASP:OD1	2.11	0.50
1:B:44:THR:OG1	1:B:46:ARG:HG3	2.11	0.50
1:A:688:PRO:HD3	1:A:694:LEU:HD11	1.92	0.50
1:D:881:ARG:HE	1:D:987:ASP:CG	2.14	0.50
1:B:464:HIS:HB2	1:B:489:GLY:HA3	1.92	0.50
1:B:416:GLU:HG3	1:B:460:ASN:O	2.12	0.50
1:B:615:PRO:O	1:B:618:THR:HG22	2.12	0.50
1:C:427:THR:HG21	1:C:462:SER:HB3	1.93	0.50
1:D:801:ILE:N	1:D:801:ILE:HD12	2.26	0.50
1:D:464:HIS:HB2	1:D:489:GLY:HA3	1.93	0.50
1:A:533:LEU:C	1:A:533:LEU:HD23	2.32	0.50
1:C:433:LEU:HB3	1:C:434:PRO:HD3	1.93	0.49
1:D:334:GLU:OE1	1:D:336:ARG:NH1	2.45	0.49
1:B:651:LEU:HD12	1:B:651:LEU:O	2.12	0.49
1:A:650:GLU:HB3	1:A:670:LEU:HD12	1.94	0.49
1:B:194:GLY:O	1:B:198:GLU:HG3	2.12	0.49
1:B:124:SER:HA	1:B:184:LEU:O	2.13	0.49
1:B:127:PHE:CE1	1:B:184:LEU:HG	2.47	0.49
1:A:651:LEU:C	1:A:651:LEU:HD12	2.32	0.49
1:A:279:ILE:HD11	1:D:422:PRO:CG	2.42	0.49
1:A:166:ARG:HG3	1:A:392:TYR:HB2	1.95	0.49
1:B:354:VAL:HG22	1:B:355:ASN:O	2.13	0.48
1:A:833:ALA:HB1	1:A:858:ILE:O	2.12	0.48
1:D:890:GLN:NE2	1:D:947:GLY:HA3	2.27	0.48
1:A:438:GLU:HG2	1:A:442:ARG:HD2	1.94	0.48
1:B:619:GLU:HA	1:B:912:ALA:HB2	1.95	0.48
1:C:630:ARG:HD3	1:C:637:GLU:OE1	2.13	0.48
1:A:433:LEU:HB3	1:A:434:PRO:HD3	1.94	0.48
1:C:738:PRO:HG2	1:C:858:ILE:O	2.14	0.48
1:A:573:GLN:HB2	1:A:602:CYS:O	2.13	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:842:TRP:HZ3	1:B:852:SER:HB3	1.79	0.48
1:C:546:LEU:HA	6:C:4132:HOH:O	2.13	0.48
1:A:887:GLN:NE2	1:A:980:GLU:O	2.44	0.48
1:A:292:ARG:C	1:A:293:LEU:HD12	2.33	0.48
1:C:166:ARG:HG3	1:C:392:TYR:CB	2.43	0.48
1:C:830:LEU:CD2	1:D:830:LEU:HD21	2.44	0.48
1:B:748:CYS:C	1:B:749:ILE:HD12	2.34	0.48
1:C:615:PRO:O	1:C:618:THR:HG22	2.14	0.48
1:D:542:ALA:HA	1:D:604:ASN:HA	1.96	0.48
1:A:755:ARG:HB2	1:A:769:TRP:HB2	1.96	0.48
1:D:895:VAL:HG21	1:D:925:MET:HG3	1.96	0.48
1:A:833:ALA:HB2	1:A:859:ASP:HA	1.96	0.48
1:C:652:LEU:HD11	1:C:698:VAL:HB	1.95	0.48
1:B:549:PHE:CE2	1:B:620:ALA:HA	2.49	0.48
1:B:636:ILE:HD13	1:B:698:VAL:HG11	1.96	0.47
1:A:830:LEU:HD22	1:B:828:ASP:HB3	1.97	0.47
1:A:619:GLU:HA	1:A:912:ALA:HB2	1.97	0.47
1:D:942:ARG:HA	1:D:953:GLY:O	2.14	0.47
1:B:143:PHE:O	1:B:168:PRO:HA	2.14	0.47
1:C:38:ASN:OD1	1:C:41:GLU:HG3	2.14	0.47
1:D:225:PHE:HA	1:D:243:GLU:O	2.14	0.47
1:B:361:PRO:HB2	1:B:576:ILE:HG12	1.95	0.47
1:A:1011:ALA:HB3	1:A:1014:TYR:CZ	2.48	0.47
1:C:828:ASP:HB3	1:D:830:LEU:HD22	1.95	0.47
1:C:549:PHE:CE2	1:C:620:ALA:HA	2.49	0.47
1:D:836:ILE:O	1:D:855:THR:HA	2.14	0.47
1:B:661:LYS:O	1:B:663:LEU:HD22	2.15	0.47
1:A:815:HIS:HE1	1:A:877:PRO:O	1.97	0.47
1:B:656:VAL:HG21	1:B:685:LEU:CD1	2.44	0.47
1:B:147:ASN:HA	1:B:148:SER:HA	1.56	0.47
1:A:570:TRP:O	1:A:607:VAL:HG22	2.14	0.47
1:B:613:PRO:HB3	1:B:617:LEU:HD23	1.96	0.47
1:A:836:ILE:O	1:A:855:THR:HA	2.15	0.47
1:D:117:GLU:HG2	6:D:4257:HOH:O	2.15	0.47
1:A:102:ASN:HD21	5:A:5130:DMS:C2	2.27	0.47
1:D:794:GLY:HA2	1:D:998:SER:O	2.15	0.47
1:C:738:PRO:HB2	1:C:834:VAL:HG23	1.97	0.47
1:D:636:ILE:HD11	1:D:682:LEU:HD11	1.97	0.47
1:C:765:LEU:HD21	1:C:768:MET:HE2	1.97	0.47
1:D:128:ASN:HB3	1:D:180:GLY:O	2.15	0.46
1:C:764:PHE:CE2	1:C:781:ARG:NH1	2.82	0.46
1:C:368:ASP:OD2	1:C:370:GLN:HB3	2.15	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:416:GLU:HA	1:C:460:ASN:O	2.15	0.46
1:A:542:ALA:HA	1:A:604:ASN:HA	1.98	0.46
1:B:13:ARG:NH1	1:C:13:ARG:HD2	2.30	0.46
1:D:301:TRP:CH2	1:D:452:SER:HA	2.49	0.46
1:C:651:LEU:HD12	1:C:668:VAL:O	2.15	0.46
1:D:476:LYS:HD2	6:D:4195:HOH:O	2.14	0.46
1:B:777:LEU:CD1	1:B:980:GLU:HG2	2.44	0.46
1:D:784:PHE:HA	1:D:881:ARG:O	2.15	0.46
1:C:595:THR:HA	1:C:596:PRO:C	2.36	0.46
1:C:573:GLN:HB2	1:C:602:CYS:O	2.15	0.46
1:B:693:GLN:OE1	1:B:724:GLU:HB2	2.15	0.46
1:D:577:LYS:O	1:D:584:PRO:HA	2.15	0.46
1:C:460:ASN:ND2	1:C:461:GLU:HG3	2.31	0.46
1:D:352:ARG:HG2	1:D:553:TRP:CH2	2.51	0.46
1:A:595:THR:HA	1:A:596:PRO:C	2.35	0.46
1:B:362:LEU:HD23	1:B:576:ILE:HB	1.97	0.46
1:A:794:GLY:HA2	1:A:998:SER:O	2.16	0.46
1:A:100:TYR:CZ	1:A:602:CYS:HB3	2.51	0.46
1:A:945:ASN:HB3	1:A:1023:LYS:HZ3	1.81	0.46
1:C:131:GLU:HG3	1:C:135:GLN:HG3	1.98	0.46
1:B:246:MET:SD	1:B:246:MET:C	2.94	0.46
1:D:144:ASP:OD2	1:D:211:ASP:HB2	2.16	0.46
1:D:26:ARG:HD2	1:D:169:SER:HA	1.98	0.46
1:D:142:ILE:HG12	1:D:170:GLU:HG2	1.97	0.46
1:B:942:ARG:HA	1:B:953:GLY:O	2.16	0.46
1:A:100:TYR:CE1	1:A:602:CYS:HB3	2.51	0.45
1:C:606:LEU:O	1:C:614:HIS:HB2	2.16	0.45
1:B:153:TRP:HB2	1:B:185:ALA:HB3	1.97	0.45
1:C:693:GLN:OE1	1:C:724:GLU:HB2	2.16	0.45
1:D:656:VAL:HG21	1:D:685:LEU:CD1	2.46	0.45
1:B:506:VAL:HG12	1:B:521:LYS:HE3	1.98	0.45
1:B:210:ARG:NH2	1:B:358:GLU:OE1	2.46	0.45
1:C:613:PRO:HB3	1:C:617:LEU:HD23	1.99	0.45
1:D:598:ASP:OD1	1:D:797:GLU:HA	2.16	0.45
1:B:646:HIS:HB3	6:B:4983:HOH:O	2.15	0.45
1:A:390:SER:HA	1:A:391:HIS:HA	1.71	0.45
1:B:246:MET:CE	1:B:250:LEU:HD23	2.46	0.45
1:C:835:LEU:HD11	1:C:855:THR:HB	1.98	0.45
1:B:546:LEU:HA	6:B:4129:HOH:O	2.16	0.45
1:A:753:ASN:O	1:A:771:GLY:N	2.48	0.45
1:C:800:ARG:HG2	1:C:800:ARG:HH11	1.81	0.45
1:A:240:LEU:C	1:A:240:LEU:HD23	2.36	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:230:ARG:NH2	1:A:241:GLU:OE2	2.49	0.45
1:D:390:SER:HA	1:D:391:HIS:HA	1.73	0.45
1:D:873:ALA:O	1:D:876:THR:HG22	2.17	0.45
1:B:801:ILE:O	1:B:803:PRO:HD3	2.17	0.45
1:A:577:LYS:O	1:A:584:PRO:HA	2.17	0.45
1:D:19:PRO:HD3	1:D:112:PRO:HB3	1.98	0.45
1:C:830:LEU:HD21	1:D:830:LEU:HD21	1.99	0.45
1:B:505:ARG:O	1:B:519:SER:HA	2.16	0.45
1:B:610:ASP:O	1:B:611:ARG:HB2	2.16	0.45
1:B:166:ARG:HG3	1:B:392:TYR:HB2	1.98	0.45
1:C:153:TRP:HB2	1:C:185:ALA:HB3	1.98	0.45
1:B:533:LEU:HD23	1:B:533:LEU:C	2.37	0.45
1:C:225:PHE:HA	1:C:243:GLU:O	2.16	0.45
1:A:797:GLU:C	1:A:799:THR:H	2.20	0.45
1:A:658:LEU:HD22	1:A:688:PRO:HB2	1.99	0.45
1:D:756:TRP:CD2	1:D:858:ILE:HD13	2.52	0.45
1:D:368:ASP:OD2	1:D:370:GLN:HB3	2.17	0.45
1:A:964:GLN:O	1:A:968:MET:HB2	2.17	0.45
1:C:965:GLN:O	1:C:969:GLU:HG3	2.16	0.45
1:C:952:ARG:NH1	1:C:952:ARG:HB2	2.32	0.44
1:B:756:TRP:CD2	1:B:858:ILE:HD13	2.51	0.44
1:A:301:TRP:CH2	1:A:452:SER:HA	2.52	0.44
1:C:568:TRP:CD2	1:C:569:ASP:HB3	2.52	0.44
1:D:196:TYR:O	1:D:417:THR:HG22	2.17	0.44
1:D:147:ASN:HA	1:D:148:SER:HA	1.62	0.44
1:B:427:THR:O	1:B:467:ASN:HB2	2.17	0.44
1:D:1011:ALA:HB3	1:D:1014:TYR:CZ	2.52	0.44
1:A:687:GLN:H	1:A:687:GLN:CD	2.20	0.44
1:A:730:LEU:HD12	1:A:730:LEU:N	2.32	0.44
1:B:1020:TRP:HD1	1:B:1021:CYS:N	2.15	0.44
1:D:577:LYS:HD3	1:D:585:TRP:CH2	2.52	0.44
1:A:587:ALA:HB1	1:A:591:ASP:CB	2.47	0.44
1:A:147:ASN:HA	1:A:148:SER:HA	1.61	0.44
1:D:755:ARG:HB2	1:D:769:TRP:HB2	2.00	0.44
1:C:653:HIS:HB3	1:C:699:ARG:NH2	2.32	0.44
1:D:688:PRO:HG3	1:D:694:LEU:HD21	2.00	0.44
1:C:143:PHE:O	1:C:168:PRO:HA	2.17	0.44
1:B:40:GLU:OE2	1:B:43:ARG:NH2	2.50	0.44
1:C:200:GLN:HA	1:C:416:GLU:OE1	2.17	0.44
1:A:370:GLN:HG3	6:A:4474:HOH:O	2.16	0.44
1:A:473:ARG:NH1	1:A:476:LYS:HB2	2.33	0.44
1:A:369:GLU:HG3	1:A:397:LEU:HD21	2.00	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:800:ARG:HH11	1:A:800:ARG:HG2	1.83	0.44
1:D:246:MET:SD	1:D:246:MET:C	2.96	0.44
1:A:422:PRO:CG	1:D:279:ILE:HD11	2.48	0.44
1:A:249:GLU:CG	1:A:251:ARG:NE	2.79	0.44
1:C:131:GLU:HG3	1:C:135:GLN:CG	2.48	0.44
1:B:546:LEU:HD22	1:B:616:ALA:HB1	1.99	0.44
1:C:505:ARG:O	1:C:519:SER:HA	2.17	0.44
1:A:920:LEU:HB3	1:A:921:PRO:HD2	2.00	0.44
1:A:770:ILE:HG23	1:A:1022:GLN:OE1	2.17	0.44
1:A:660:GLY:O	1:A:662:PRO:HD3	2.18	0.44
1:D:610:ASP:O	1:D:611:ARG:HB2	2.18	0.44
1:A:945:ASN:OD1	1:A:1023:LYS:HD3	2.18	0.43
1:A:802:ASP:O	1:A:808:GLU:HG3	2.18	0.43
1:C:326:GLU:OE1	1:C:326:GLU:HA	2.18	0.43
1:D:390:SER:HB2	1:D:391:HIS:CE1	2.54	0.43
1:B:91:GLN:HG3	1:B:96:ASP:OD1	2.17	0.43
1:A:842:TRP:HZ3	1:A:852:SER:HB3	1.82	0.43
1:C:525:SER:O	1:D:561:ARG:HD3	2.18	0.43
1:A:610:ASP:O	1:A:611:ARG:HB2	2.18	0.43
1:B:824:GLN:HB3	1:B:839:ALA:HB3	2.00	0.43
1:D:240:LEU:HD23	1:D:240:LEU:C	2.38	0.43
1:D:355:ASN:OD1	1:D:388:ARG:HD3	2.18	0.43
1:C:777:LEU:HB2	1:C:887:GLN:HG2	2.00	0.43
1:B:147:ASN:HB3	1:B:206:SER:HA	2.01	0.43
1:B:833:ALA:HB1	1:B:858:ILE:O	2.19	0.43
1:D:673:ALA:HB1	1:D:674:PRO:HD2	2.00	0.43
1:D:687:GLN:H	1:D:687:GLN:CD	2.22	0.43
1:B:598:ASP:OD1	1:B:797:GLU:HA	2.17	0.43
1:D:568:TRP:HA	1:D:569:ASP:HA	1.78	0.43
1:D:843:GLN:HA	1:D:847:LYS:O	2.18	0.43
1:C:59:ARG:HB2	1:C:124:SER:OG	2.19	0.43
1:D:524:LEU:HD11	1:D:562:LEU:HG	2.01	0.43
1:C:777:LEU:HD11	1:C:980:GLU:HG2	2.01	0.43
1:A:634:GLN:HE21	1:A:684:GLU:HA	1.82	0.43
1:B:937:LEU:O	1:B:938:ARG:HD2	2.17	0.43
1:A:91:GLN:HG3	1:A:96:ASP:OD1	2.19	0.43
1:B:19:PRO:HD3	1:B:112:PRO:CB	2.48	0.43
1:A:427:THR:O	1:A:467:ASN:HB2	2.19	0.43
1:B:352:ARG:HG2	1:B:553:TRP:CH2	2.54	0.43
1:A:622:HIS:ND1	1:A:625:GLN:OE1	2.49	0.43
1:D:427:THR:O	1:D:467:ASN:HB2	2.18	0.43
1:C:245:GLN:HG2	1:C:288:ARG:CG	2.48	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:568:TRP:CD2	1:D:569:ASP:HB3	2.53	0.42
1:B:788:PRO:HD2	1:B:968:MET:HG3	2.01	0.42
1:B:657:ALA:HA	1:B:663:LEU:HD22	2.00	0.42
1:C:749:ILE:N	1:C:749:ILE:CD1	2.82	0.42
1:D:890:GLN:HB2	6:D:4659:HOH:O	2.18	0.42
1:A:256:VAL:HA	1:A:314:GLU:O	2.19	0.42
1:A:128:ASN:HA	1:A:180:GLY:O	2.19	0.42
1:D:100:TYR:CE1	1:D:602:CYS:HB3	2.54	0.42
1:D:52:ARG:O	1:D:213:SER:HB2	2.19	0.42
1:B:638:VAL:O	1:B:677:LYS:HA	2.19	0.42
1:A:173:LEU:O	1:A:177:LEU:HG	2.20	0.42
1:D:619:GLU:HA	1:D:912:ALA:HB2	2.01	0.42
1:A:361:PRO:HB3	1:A:609:ALA:HB1	2.02	0.42
1:D:984:LEU:HD21	1:D:986:ILE:HD11	2.01	0.42
1:A:225:PHE:HA	1:A:243:GLU:O	2.20	0.42
1:D:533:LEU:HD23	1:D:533:LEU:C	2.40	0.42
1:D:340:GLY:O	1:D:561:ARG:HG2	2.20	0.42
1:A:16:TRP:CG	1:A:189:LEU:HD13	2.55	0.42
1:A:87:PRO:HA	1:A:208:ILE:O	2.20	0.42
1:A:618:THR:HG22	1:A:912:ALA:HB1	2.01	0.42
1:A:703:PRO:O	1:A:711:ALA:HB1	2.20	0.42
1:B:944:LEU:O	1:B:950:GLN:HA	2.20	0.42
1:B:896:ASN:HB3	1:B:945:ASN:HB2	2.01	0.42
1:C:661:LYS:HA	1:C:662:PRO:HD3	1.92	0.42
1:C:890:GLN:HE22	1:C:948:PRO:HD3	1.84	0.42
1:A:598:ASP:OD1	1:A:797:GLU:HA	2.20	0.41
1:B:815:HIS:HE1	1:B:877:PRO:O	2.02	0.41
1:A:738:PRO:HD2	1:A:833:ALA:HA	2.03	0.41
1:A:433:LEU:N	1:A:434:PRO:CD	2.83	0.41
1:A:773:LYS:HG2	1:A:774:LYS:O	2.20	0.41
1:B:581:ASN:HB2	1:B:583:ASN:ND2	2.35	0.41
1:C:105:TYR:CE1	1:C:199:ASP:HB2	2.54	0.41
1:C:200:GLN:HG3	1:C:416:GLU:OE1	2.20	0.41
1:B:658:LEU:HD22	1:B:688:PRO:HB2	2.02	0.41
1:B:506:VAL:CG1	1:B:521:LYS:HE3	2.50	0.41
1:B:634:GLN:HB2	1:B:681:GLU:OE2	2.21	0.41
1:C:19:PRO:HD3	1:C:112:PRO:CB	2.50	0.41
1:B:105:TYR:CE1	1:B:199:ASP:HB2	2.56	0.41
1:A:754:LYS:HA	1:A:769:TRP:O	2.19	0.41
1:A:656:VAL:HG21	1:A:685:LEU:CD1	2.50	0.41
1:B:128:ASN:HA	1:B:180:GLY:O	2.21	0.41
1:D:738:PRO:HG3	1:D:751:LEU:HD13	2.01	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:784:PHE:HA	1:C:881:ARG:O	2.21	0.41
1:B:390:SER:HA	1:B:391:HIS:HA	1.81	0.41
1:A:525:SER:O	1:B:561:ARG:HD3	2.21	0.41
1:C:246:MET:C	1:C:246:MET:SD	2.99	0.41
1:C:952:ARG:HH11	1:C:952:ARG:CB	2.33	0.41
1:D:166:ARG:HG3	1:D:392:TYR:HB2	2.02	0.41
1:C:482:ARG:HA	1:C:483:PRO:HD3	1.97	0.41
1:D:458:LEU:HD11	1:D:472:TYR:HB2	2.03	0.41
1:D:863:GLN:NE2	1:D:952:ARG:NH2	2.67	0.41
1:B:674:PRO:O	1:B:675:GLN:HB2	2.21	0.41
1:A:390:SER:HB2	1:A:391:HIS:CE1	2.56	0.41
1:B:895:VAL:HG21	1:B:925:MET:HG3	2.03	0.41
1:A:568:TRP:CD2	1:A:569:ASP:HB3	2.56	0.41
1:B:223:SER:O	1:B:224:ASP:HB2	2.21	0.41
1:A:292:ARG:HH12	5:A:5012:DMS:C1	2.34	0.41
1:C:100:TYR:CE1	1:C:602:CYS:HB3	2.56	0.41
1:A:587:ALA:HB1	1:A:591:ASP:HB2	2.03	0.41
1:B:19:PRO:HD3	1:B:112:PRO:HB3	2.02	0.41
1:B:141:ILE:HB	1:B:173:LEU:HD11	2.03	0.41
1:C:638:VAL:O	1:C:677:LYS:HA	2.21	0.41
6:A:4541:HOH:O	1:D:463:GLY:HA2	2.20	0.41
1:D:482:ARG:HA	1:D:483:PRO:HD3	1.96	0.41
1:A:335:VAL:HG22	1:A:344:LEU:HD12	2.03	0.41
1:A:126:THR:HA	1:A:182:ASN:O	2.21	0.41
1:D:80:GLU:CD	1:D:80:GLU:H	2.25	0.41
1:A:793:ILE:HG22	1:A:795:VAL:HG22	2.03	0.41
1:A:701:VAL:O	1:A:703:PRO:HD3	2.20	0.41
1:A:890:GLN:HG2	1:A:891:VAL:N	2.35	0.41
1:C:749:ILE:HD11	1:C:836:ILE:HD11	2.03	0.40
1:A:651:LEU:O	1:A:651:LEU:HD12	2.21	0.40
1:C:794:GLY:HA2	1:C:998:SER:O	2.21	0.40
1:C:802:ASP:O	1:C:808:GLU:HG3	2.21	0.40
1:C:542:ALA:HA	1:C:604:ASN:HA	2.03	0.40
1:A:499:ILE:HG22	1:A:501:PRO:HD3	2.03	0.40
1:B:658:LEU:N	1:B:663:LEU:CD2	2.84	0.40
1:B:251:ARG:HB2	1:B:254:LEU:HG	2.03	0.40
1:B:163:GLN:O	1:B:164:ASP:HB3	2.21	0.40
1:A:900:LEU:CD1	1:A:910:LEU:HD22	2.51	0.40
1:A:896:ASN:HD21	1:A:917:ARG:HD2	1.86	0.40
1:D:429:ASP:OD1	1:D:431:ARG:HG3	2.21	0.40
1:D:742:THR:HG22	1:D:743:SER:N	2.37	0.40
1:C:952:ARG:HH11	1:C:952:ARG:HB2	1.87	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:847:LYS:HE2	1:B:848:THR:O	2.21	0.40
1:C:619:GLU:HA	1:C:912:ALA:HB2	2.02	0.40
1:A:768:MET:HE3	1:A:1020:TRP:CZ2	2.56	0.40
1:D:773:LYS:HA	1:D:773:LYS:HD2	1.94	0.40
1:A:291:LEU:HD22	1:A:291:LEU:N	2.37	0.40
1:A:88:SER:HA	1:A:366:VAL:CG2	2.48	0.40
1:C:533:LEU:HD23	1:C:534:ILE:N	2.36	0.40
1:A:533:LEU:HD23	1:A:534:ILE:N	2.36	0.40
1:C:546:LEU:HD22	1:C:616:ALA:HB1	2.04	0.40
1:A:87:PRO:HB2	1:A:209:PHE:C	2.42	0.40
1:A:59:ARG:HB2	1:A:124:SER:OG	2.21	0.40
1:A:742:THR:HG22	1:A:743:SER:N	2.36	0.40
1:B:372:MET:CE	1:B:397:LEU:HD23	2.52	0.40
1:B:789:LEU:HD11	1:B:993:ILE:HG22	2.03	0.40
6:B:4559:HOH:O	1:C:463:GLY:HA2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1009/1023 (99%)	963 (95%)	45 (4%)	1 (0%)	59	66
1	B	1009/1023 (99%)	961 (95%)	48 (5%)	0	100	100
1	C	1009/1023 (99%)	962 (95%)	47 (5%)	0	100	100
1	D	1009/1023 (99%)	965 (96%)	43 (4%)	1 (0%)	59	66
All	All	4036/4092 (99%)	3851 (95%)	183 (4%)	2 (0%)	100	100

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	798	ALA
1	D	211	ASP

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	863/874 (99%)	858 (99%)	5 (1%)	92	97
1	B	863/874 (99%)	858 (99%)	5 (1%)	92	97
1	C	863/874 (99%)	857 (99%)	6 (1%)	91	96
1	D	863/874 (99%)	857 (99%)	6 (1%)	91	96
All	All	3452/3496 (99%)	3430 (99%)	22 (1%)	92	97

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	277	GLU
1	A	333	ARG
1	A	519	SER
1	A	546	LEU
1	A	910	LEU
1	B	251	ARG
1	B	333	ARG
1	B	546	LEU
1	B	604	ASN
1	B	956	GLN
1	C	112	PRO
1	C	333	ARG
1	C	344	LEU
1	C	546	LEU
1	C	773	LYS
1	C	804	ASN
1	D	76	CYS
1	D	333	ARG
1	D	546	LEU
1	D	655	MET
1	D	663	LEU
1	D	885	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	102	ASN
1	A	634	GLN
1	A	757	GLN
1	A	1017	GLN
1	B	262	GLN
1	B	583	ASN
1	B	804	ASN
1	B	965	GLN
1	C	817	GLN
1	C	863	GLN
1	D	675	GLN
1	D	863	GLN

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 ⓘ

Of 157 ligands modelled in this entry, 27 are monoatomic - leaving 130 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	149	A	2001	4	12,12,12	1.33	2 (16%)	17,17,17	0.84	0
5	DMS	A	5001	-	3,3,3	0.20	0	3,3,3	0.56	0
5	DMS	A	5002	-	3,3,3	0.18	0	3,3,3	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	DMS	A	5003	-	3,3,3	0.20	0	3,3,3	0.54	0
5	DMS	A	5004	-	3,3,3	0.22	0	3,3,3	0.57	0
5	DMS	A	5005	-	3,3,3	0.23	0	3,3,3	0.52	0
5	DMS	A	5006	-	3,3,3	0.23	0	3,3,3	0.54	0
5	DMS	A	5007	-	3,3,3	0.23	0	3,3,3	0.53	0
5	DMS	A	5008	-	3,3,3	0.23	0	3,3,3	0.57	0
5	DMS	A	5009	-	3,3,3	0.21	0	3,3,3	0.55	0
5	DMS	A	5010	-	3,3,3	0.23	0	3,3,3	0.55	0
5	DMS	A	5011	-	3,3,3	0.24	0	3,3,3	0.53	0
5	DMS	A	5012	-	3,3,3	0.22	0	3,3,3	0.55	0
5	DMS	A	5014	-	3,3,3	0.24	0	3,3,3	0.56	0
5	DMS	A	5015	-	3,3,3	0.22	0	3,3,3	0.56	0
5	DMS	A	5016	-	3,3,3	0.24	0	3,3,3	0.55	0
5	DMS	A	5017	-	3,3,3	0.22	0	3,3,3	0.58	0
5	DMS	A	5018	4	3,3,3	0.23	0	3,3,3	0.57	0
5	DMS	A	5019	-	3,3,3	0.21	0	3,3,3	0.55	0
5	DMS	A	5020	-	3,3,3	0.18	0	3,3,3	0.55	0
5	DMS	A	5022	-	3,3,3	0.24	0	3,3,3	0.56	0
5	DMS	A	5023	-	3,3,3	0.21	0	3,3,3	0.53	0
5	DMS	A	5024	-	3,3,3	0.24	0	3,3,3	0.57	0
5	DMS	A	5125	-	3,3,3	0.23	0	3,3,3	0.55	0
5	DMS	A	5126	-	3,3,3	0.22	0	3,3,3	0.56	0
5	DMS	A	5127	-	3,3,3	0.22	0	3,3,3	0.57	0
5	DMS	A	5128	-	3,3,3	0.25	0	3,3,3	0.58	0
5	DMS	A	5129	-	3,3,3	0.24	0	3,3,3	0.56	0
5	DMS	A	5130	-	3,3,3	0.14	0	3,3,3	0.53	0
5	DMS	A	5231	-	3,3,3	0.20	0	3,3,3	0.56	0
5	DMS	A	5232	-	3,3,3	0.23	0	3,3,3	0.59	0
2	149	B	2001	4	12,12,12	1.17	2 (16%)	17,17,17	0.85	0
5	DMS	B	5001	-	3,3,3	0.24	0	3,3,3	0.54	0
5	DMS	B	5002	-	3,3,3	0.19	0	3,3,3	0.52	0
5	DMS	B	5003	-	3,3,3	0.22	0	3,3,3	0.57	0
5	DMS	B	5004	-	3,3,3	0.18	0	3,3,3	0.53	0
5	DMS	B	5005	-	3,3,3	0.21	0	3,3,3	0.54	0
5	DMS	B	5006	-	3,3,3	0.26	0	3,3,3	0.54	0
5	DMS	B	5007	-	3,3,3	0.21	0	3,3,3	0.54	0
5	DMS	B	5008	-	3,3,3	0.18	0	3,3,3	0.55	0
5	DMS	B	5009	-	3,3,3	0.23	0	3,3,3	0.56	0
5	DMS	B	5010	-	3,3,3	0.23	0	3,3,3	0.56	0
5	DMS	B	5011	-	3,3,3	0.21	0	3,3,3	0.52	0
5	DMS	B	5012	-	3,3,3	0.25	0	3,3,3	0.54	0
5	DMS	B	5013	-	3,3,3	0.23	0	3,3,3	0.56	0
5	DMS	B	5014	-	3,3,3	0.20	0	3,3,3	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	DMS	B	5015	-	3,3,3	0.22	0	3,3,3	0.55	0
5	DMS	B	5016	-	3,3,3	0.22	0	3,3,3	0.54	0
5	DMS	B	5017	-	3,3,3	0.23	0	3,3,3	0.57	0
5	DMS	B	5018	-	3,3,3	0.21	0	3,3,3	0.55	0
5	DMS	B	5019	-	3,3,3	0.22	0	3,3,3	0.55	0
5	DMS	B	5020	-	3,3,3	0.22	0	3,3,3	0.55	0
5	DMS	B	5021	4	3,3,3	0.23	0	3,3,3	0.55	0
5	DMS	B	5022	-	3,3,3	0.23	0	3,3,3	0.54	0
5	DMS	B	5023	-	3,3,3	0.22	0	3,3,3	0.56	0
5	DMS	B	5024	-	3,3,3	0.21	0	3,3,3	0.55	0
5	DMS	B	5025	-	3,3,3	0.24	0	3,3,3	0.58	0
5	DMS	B	5026	-	3,3,3	0.22	0	3,3,3	0.56	0
5	DMS	B	5027	-	3,3,3	0.24	0	3,3,3	0.55	0
5	DMS	B	5028	-	3,3,3	0.23	0	3,3,3	0.56	0
5	DMS	B	5029	-	3,3,3	0.23	0	3,3,3	0.55	0
5	DMS	B	5034	-	3,3,3	0.22	0	3,3,3	0.54	0
5	DMS	B	5131	-	3,3,3	0.20	0	3,3,3	0.54	0
5	DMS	B	5132	-	3,3,3	0.18	0	3,3,3	0.54	0
5	DMS	B	5133	-	3,3,3	0.22	0	3,3,3	0.55	0
5	DMS	C	1024	-	3,3,3	0.23	0	3,3,3	0.55	0
2	149	C	2001	4	12,12,12	1.44	3 (25%)	17,17,17	0.82	0
5	DMS	C	5001	-	3,3,3	0.21	0	3,3,3	0.54	0
5	DMS	C	5002	-	3,3,3	0.18	0	3,3,3	0.54	0
5	DMS	C	5003	-	3,3,3	0.19	0	3,3,3	0.55	0
5	DMS	C	5004	-	3,3,3	0.21	0	3,3,3	0.54	0
5	DMS	C	5005	-	3,3,3	0.21	0	3,3,3	0.53	0
5	DMS	C	5006	-	3,3,3	0.24	0	3,3,3	0.57	0
5	DMS	C	5007	-	3,3,3	0.23	0	3,3,3	0.54	0
5	DMS	C	5008	-	3,3,3	0.22	0	3,3,3	0.54	0
5	DMS	C	5009	-	3,3,3	0.23	0	3,3,3	0.55	0
5	DMS	C	5010	-	3,3,3	0.22	0	3,3,3	0.54	0
5	DMS	C	5011	-	3,3,3	0.22	0	3,3,3	0.54	0
5	DMS	C	5012	-	3,3,3	0.22	0	3,3,3	0.55	0
5	DMS	C	5013	-	3,3,3	0.24	0	3,3,3	0.54	0
5	DMS	C	5014	-	3,3,3	0.22	0	3,3,3	0.55	0
5	DMS	C	5015	-	3,3,3	0.21	0	3,3,3	0.55	0
5	DMS	C	5016	-	3,3,3	0.23	0	3,3,3	0.56	0
5	DMS	C	5017	-	3,3,3	0.26	0	3,3,3	0.56	0
5	DMS	C	5018	-	3,3,3	0.20	0	3,3,3	0.56	0
5	DMS	C	5019	-	3,3,3	0.23	0	3,3,3	0.55	0
5	DMS	C	5020	-	3,3,3	0.22	0	3,3,3	0.55	0
5	DMS	C	5021	4	3,3,3	0.22	0	3,3,3	0.56	0
5	DMS	C	5022	-	3,3,3	0.20	0	3,3,3	0.53	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	DMS	C	5023	-	3,3,3	0.20	0	3,3,3	0.56	0
5	DMS	C	5024	-	3,3,3	0.21	0	3,3,3	0.57	0
5	DMS	C	5025	-	3,3,3	0.23	0	3,3,3	0.56	0
5	DMS	C	5027	-	3,3,3	0.25	0	3,3,3	0.55	0
5	DMS	C	5028	-	3,3,3	0.23	0	3,3,3	0.55	0
5	DMS	C	5030	-	3,3,3	0.26	0	3,3,3	0.57	0
5	DMS	C	5031	-	3,3,3	0.21	0	3,3,3	0.55	0
5	DMS	C	5032	-	3,3,3	0.20	0	3,3,3	0.55	0
5	DMS	C	5033	-	3,3,3	0.25	0	3,3,3	0.56	0
5	DMS	C	5134	-	3,3,3	0.25	0	3,3,3	0.56	0
2	149	D	2001	4	12,12,12	1.16	1 (8%)	17,17,17	0.86	0
5	DMS	D	5001	-	3,3,3	0.22	0	3,3,3	0.59	0
5	DMS	D	5002	-	3,3,3	0.17	0	3,3,3	0.48	0
5	DMS	D	5003	-	3,3,3	0.20	0	3,3,3	0.55	0
5	DMS	D	5004	-	3,3,3	0.17	0	3,3,3	0.55	0
5	DMS	D	5005	-	3,3,3	0.24	0	3,3,3	0.53	0
5	DMS	D	5006	-	3,3,3	0.23	0	3,3,3	0.53	0
5	DMS	D	5007	-	3,3,3	0.20	0	3,3,3	0.54	0
5	DMS	D	5008	-	3,3,3	0.22	0	3,3,3	0.56	0
5	DMS	D	5009	-	3,3,3	0.20	0	3,3,3	0.55	0
5	DMS	D	5010	-	3,3,3	0.23	0	3,3,3	0.55	0
5	DMS	D	5011	-	3,3,3	0.23	0	3,3,3	0.54	0
5	DMS	D	5012	-	3,3,3	0.23	0	3,3,3	0.55	0
5	DMS	D	5013	-	3,3,3	0.21	0	3,3,3	0.55	0
5	DMS	D	5014	-	3,3,3	0.22	0	3,3,3	0.55	0
5	DMS	D	5015	-	3,3,3	0.21	0	3,3,3	0.56	0
5	DMS	D	5016	-	3,3,3	0.21	0	3,3,3	0.54	0
5	DMS	D	5017	-	3,3,3	0.23	0	3,3,3	0.56	0
5	DMS	D	5018	-	3,3,3	0.24	0	3,3,3	0.57	0
5	DMS	D	5019	-	3,3,3	0.24	0	3,3,3	0.55	0
5	DMS	D	5020	-	3,3,3	0.22	0	3,3,3	0.54	0
5	DMS	D	5021	-	3,3,3	0.20	0	3,3,3	0.54	0
5	DMS	D	5022	-	3,3,3	0.22	0	3,3,3	0.56	0
5	DMS	D	5023	-	3,3,3	0.19	0	3,3,3	0.54	0
5	DMS	D	5024	-	3,3,3	0.22	0	3,3,3	0.57	0
5	DMS	D	5025	-	3,3,3	0.24	0	3,3,3	0.55	0
5	DMS	D	5026	-	3,3,3	0.21	0	3,3,3	0.55	0
5	DMS	D	5027	-	3,3,3	0.23	0	3,3,3	0.55	0
5	DMS	D	5028	-	3,3,3	0.23	0	3,3,3	0.57	0
5	DMS	D	5229	-	3,3,3	0.24	0	3,3,3	0.56	0
5	DMS	D	5230	-	3,3,3	0.21	0	3,3,3	0.57	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	149	A	2001	4	-	0/2/22/22	0/1/1/1
5	DMS	A	5001	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5002	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5003	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5004	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5005	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5006	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5007	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5008	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5009	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5010	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5011	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5012	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5014	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5015	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5016	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5017	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5018	4	-	0/0/0/0	0/0/0/0
5	DMS	A	5019	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5020	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5022	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5023	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5024	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5125	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5126	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5127	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5128	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5129	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5130	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5231	-	-	0/0/0/0	0/0/0/0
5	DMS	A	5232	-	-	0/0/0/0	0/0/0/0
2	149	B	2001	4	-	0/2/22/22	0/1/1/1
5	DMS	B	5001	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5002	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5003	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5004	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5005	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5006	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5007	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5008	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	DMS	B	5009	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5010	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5011	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5012	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5013	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5014	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5015	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5016	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5017	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5018	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5019	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5020	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5021	4	-	0/0/0/0	0/0/0/0
5	DMS	B	5022	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5023	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5024	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5025	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5026	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5027	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5028	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5029	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5034	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5131	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5132	-	-	0/0/0/0	0/0/0/0
5	DMS	B	5133	-	-	0/0/0/0	0/0/0/0
5	DMS	C	1024	-	-	0/0/0/0	0/0/0/0
2	149	C	2001	4	-	0/2/22/22	0/1/1/1
5	DMS	C	5001	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5002	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5003	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5004	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5005	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5006	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5007	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5008	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5009	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5010	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5011	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5012	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5013	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5014	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5015	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	DMS	C	5016	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5017	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5018	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5019	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5020	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5021	4	-	0/0/0/0	0/0/0/0
5	DMS	C	5022	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5023	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5024	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5025	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5027	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5028	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5030	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5031	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5032	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5033	-	-	0/0/0/0	0/0/0/0
5	DMS	C	5134	-	-	0/0/0/0	0/0/0/0
2	149	D	2001	4	-	0/2/22/22	0/1/1/1
5	DMS	D	5001	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5002	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5003	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5004	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5005	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5006	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5007	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5008	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5009	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5010	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5011	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5012	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5013	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5014	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5015	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5016	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5017	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5018	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5019	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5020	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5021	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5022	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5023	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5024	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	DMS	D	5025	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5026	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5027	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5028	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5229	-	-	0/0/0/0	0/0/0/0
5	DMS	D	5230	-	-	0/0/0/0	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2001	149	C2-C1	3.11	1.54	1.51
2	A	2001	149	O5-C1	2.92	1.39	1.34
2	C	2001	149	O5-C1	2.80	1.38	1.34
2	B	2001	149	O5-C1	2.72	1.38	1.34
2	D	2001	149	O5-C1	2.72	1.38	1.34
2	A	2001	149	C2-C1	2.59	1.54	1.51
2	C	2001	149	O5-C5	2.09	1.49	1.46
2	B	2001	149	C2-C1	2.01	1.53	1.51

There are no bond angle outliers.

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	1011/1023 (98%)	-0.35	26 (2%) 53 53	7, 20, 41, 66	0
1	B	1011/1023 (98%)	-0.42	22 (2%) 59 59	5, 18, 39, 66	0
1	C	1011/1023 (98%)	-0.47	21 (2%) 60 61	6, 17, 37, 73	0
1	D	1011/1023 (98%)	-0.45	22 (2%) 59 59	6, 18, 38, 66	0
All	All	4044/4092 (98%)	-0.42	91 (2%) 59 58	5, 18, 39, 73	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	732	ALA	6.1
1	A	735	HIS	5.9
1	C	732	ALA	5.8
1	B	732	ALA	5.6
1	A	732	ALA	5.4
1	C	801	ILE	5.3
1	C	733	ALA	5.2
1	A	795	VAL	5.2
1	D	689	GLU	5.0
1	C	735	HIS	5.0
1	C	800	ARG	4.9
1	D	801	ILE	4.8
1	B	731	PRO	4.8
1	A	730	LEU	4.7
1	A	799	THR	4.7
1	A	801	ILE	4.7
1	A	1023	LYS	4.6
1	D	731	PRO	4.6
1	C	799	THR	4.5
1	C	689	GLU	4.5
1	C	731	PRO	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	800	ARG	4.4
1	A	734	SER	4.4
1	B	733	ALA	4.2
1	B	686	PRO	4.1
1	D	735	HIS	4.0
1	A	689	GLU	3.9
1	D	730	LEU	3.9
1	A	687	GLN	3.8
1	B	689	GLU	3.8
1	B	687	GLN	3.7
1	A	684	GLU	3.6
1	A	733	ALA	3.6
1	B	800	ARG	3.5
1	D	800	ARG	3.5
1	B	684	GLU	3.5
1	C	580	GLU	3.5
1	D	687	GLN	3.4
1	B	580	GLU	3.4
1	A	686	PRO	3.4
1	C	730	LEU	3.3
1	A	731	PRO	3.3
1	D	734	SER	3.3
1	D	684	GLU	3.3
1	C	744	GLU	3.3
1	D	772	ASP	3.3
1	C	772	ASP	3.2
1	C	734	SER	3.2
1	D	580	GLU	3.1
1	C	687	GLN	3.0
1	D	1023	LYS	3.0
1	C	745	MET	3.0
1	B	730	LEU	3.0
1	B	1023	LYS	3.0
1	A	580	GLU	3.0
1	A	729	THR	2.9
1	B	734	SER	2.9
1	B	801	ILE	2.8
1	A	771	GLY	2.8
1	B	690	SER	2.8
1	D	733	ALA	2.8
1	A	71	GLU	2.7
1	C	1022	GLN	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	1022	GLN	2.6
1	D	799	THR	2.6
1	D	831	ALA	2.6
1	D	686	PRO	2.5
1	B	745	MET	2.5
1	A	76	CYS	2.5
1	D	845	GLN	2.5
1	B	735	HIS	2.5
1	B	772	ASP	2.5
1	D	1022	GLN	2.4
1	A	819	GLU	2.4
1	B	799	THR	2.4
1	C	684	GLU	2.4
1	B	795	VAL	2.3
1	C	817	GLN	2.3
1	C	795	VAL	2.3
1	A	49	GLN	2.3
1	B	1022	GLN	2.3
1	B	688	PRO	2.2
1	A	772	ASP	2.2
1	C	686	PRO	2.1
1	C	1023	LYS	2.1
1	D	795	VAL	2.1
1	B	744	GLU	2.1
1	A	178	ARG	2.1
1	D	71	GLU	2.1
1	A	249	GLU	2.0
1	D	634	GLN	2.0

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

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

6.3 Carbohydrates ⓘ

There are no 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
5	DMS	A	5010	4/4	0.30	18.12	71,71,71,72	0
5	DMS	D	5013	4/4	0.34	15.21	87,87,87,88	0
5	DMS	B	5026	4/4	0.21	14.36	55,55,55,56	0
5	DMS	D	5015	4/4	0.31	14.29	78,79,79,79	0
5	DMS	C	5009	4/4	0.24	13.05	90,90,90,91	0
5	DMS	D	5230	4/4	0.18	13.00	54,54,54,55	0
5	DMS	B	5015	4/4	0.18	11.98	59,59,60,60	0
5	DMS	B	5022	4/4	0.26	11.72	89,89,89,90	0
5	DMS	A	5024	4/4	0.27	11.44	79,79,79,79	0
5	DMS	C	5019	4/4	0.18	10.92	51,52,52,52	0
5	DMS	D	5010	4/4	0.24	10.45	71,71,71,72	0
5	DMS	B	5019	4/4	0.17	10.11	60,60,60,61	0
5	DMS	D	5007	4/4	0.31	9.86	86,87,87,87	0
5	DMS	B	5021	4/4	0.32	9.73	66,66,66,67	0
5	DMS	C	5014	4/4	0.25	9.44	65,65,66,66	0
5	DMS	C	5032	4/4	0.20	8.93	61,62,62,63	0
5	DMS	C	5022	4/4	0.20	8.36	65,65,65,66	0
5	DMS	A	5019	4/4	0.26	8.05	84,84,84,84	0
5	DMS	D	5024	4/4	0.29	8.03	85,85,86,86	0
5	DMS	B	5020	4/4	0.25	7.80	67,68,68,68	0
5	DMS	A	5016	4/4	0.22	7.78	57,57,57,58	0
5	DMS	C	5033	4/4	0.16	7.73	58,58,59,60	0
5	DMS	A	5017	4/4	0.24	7.66	59,59,59,60	0
5	DMS	C	5012	4/4	0.35	7.53	97,97,97,97	0
5	DMS	C	1024	4/4	0.21	7.45	63,64,64,65	0
5	DMS	A	5006	4/4	0.15	7.32	57,58,58,58	0
5	DMS	C	5017	4/4	0.16	7.30	65,65,66,67	0
5	DMS	D	5021	4/4	0.20	7.12	83,83,83,83	0
5	DMS	A	5125	4/4	0.28	7.09	72,72,72,73	0
5	DMS	C	5016	4/4	0.27	6.46	68,68,69,69	0
5	DMS	D	5019	4/4	0.20	6.29	56,57,57,57	0
5	DMS	C	5013	4/4	0.18	6.23	38,40,40,40	0
5	DMS	A	5007	4/4	0.22	6.14	65,66,66,66	0
5	DMS	B	5007	4/4	0.22	6.05	75,75,76,76	0
5	DMS	A	5002	4/4	0.17	6.03	45,46,46,47	0
5	DMS	C	5006	4/4	0.17	5.86	56,57,58,58	0
5	DMS	C	5018	4/4	0.19	5.65	66,67,67,67	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	DMS	B	5025	4/4	0.21	5.53	70,70,70,72	0
5	DMS	B	5009	4/4	0.18	5.38	39,40,41,41	0
5	DMS	C	5028	4/4	0.23	5.34	84,85,85,85	0
5	DMS	B	5006	4/4	0.23	5.29	63,63,64,64	0
5	DMS	B	5003	4/4	0.14	5.09	39,40,41,41	0
5	DMS	C	5021	4/4	0.22	4.74	60,60,60,61	0
5	DMS	B	5131	4/4	0.15	4.60	58,58,58,59	0
5	DMS	D	5004	4/4	0.17	4.55	52,52,53,54	0
5	DMS	B	5017	4/4	0.30	4.48	74,74,74,75	0
5	DMS	A	5014	4/4	0.19	4.44	49,51,51,51	0
5	DMS	C	5004	4/4	0.15	4.38	36,37,39,39	0
5	DMS	C	5020	4/4	0.19	4.23	68,68,68,69	0
5	DMS	C	5003	4/4	0.13	4.17	36,37,38,39	0
5	DMS	D	5014	4/4	0.20	4.11	58,59,60,60	0
4	NA	D	3104	1/1	0.18	4.10	29,29,29,29	0
5	DMS	D	5229	4/4	0.21	3.97	67,67,69,69	0
5	DMS	A	5127	4/4	0.26	3.62	51,53,53,53	0
5	DMS	A	5009	4/4	0.19	3.60	44,45,46,47	0
5	DMS	B	5004	4/4	0.14	3.59	51,52,53,53	0
5	DMS	B	5018	4/4	0.15	3.52	45,46,46,48	0
5	DMS	A	5004	4/4	0.17	3.50	57,57,58,59	0
5	DMS	D	5026	4/4	0.23	3.24	113,113,113,113	0
5	DMS	D	5020	4/4	0.15	3.10	65,66,66,66	0
5	DMS	D	5023	4/4	0.30	3.05	81,82,82,82	0
5	DMS	A	5015	4/4	0.18	2.96	65,65,65,66	0
5	DMS	D	5018	4/4	0.18	2.94	71,71,71,72	0
2	149	A	2001	12/12	0.11	2.74	11,13,18,18	0
5	DMS	B	5132	4/4	0.14	2.73	58,58,59,59	0
5	DMS	B	5011	4/4	0.13	2.67	28,28,29,29	0
5	DMS	B	5002	4/4	0.12	2.67	38,38,40,41	0
5	DMS	D	5027	4/4	0.25	2.65	75,75,76,76	0
5	DMS	B	5013	4/4	0.31	2.57	109,109,109,109	0
5	DMS	D	5022	4/4	0.16	2.52	62,62,62,63	0
5	DMS	C	5008	4/4	0.15	2.50	55,55,55,55	0
5	DMS	D	5002	4/4	0.13	2.44	30,31,32,32	0
5	DMS	C	5027	4/4	0.20	2.43	65,66,66,66	0
5	DMS	C	5134	4/4	0.20	2.27	40,41,41,43	0
5	DMS	B	5028	4/4	0.19	2.25	53,53,54,55	0
2	149	D	2001	12/12	0.10	2.18	8,13,15,20	0
5	DMS	B	5027	4/4	0.18	2.05	52,52,52,52	0
5	DMS	B	5014	4/4	0.15	1.89	40,40,41,42	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	DMS	A	5022	4/4	0.25	1.82	65,65,66,66	0
3	MG	B	3003	1/1	0.12	1.74	59,59,59,59	0
5	DMS	C	5002	4/4	0.11	1.69	32,33,34,35	0
5	DMS	C	5030	4/4	0.17	1.58	62,63,63,64	0
5	DMS	C	5024	4/4	0.23	1.56	66,66,66,67	0
5	DMS	C	5031	4/4	0.23	1.55	69,69,70,70	0
5	DMS	D	5009	4/4	0.16	1.50	44,45,45,46	0
5	DMS	B	5023	4/4	0.15	1.46	41,42,42,43	0
5	DMS	B	5010	4/4	0.15	1.43	72,72,73,73	0
5	DMS	A	5128	4/4	0.24	1.35	54,54,55,55	0
5	DMS	C	5007	4/4	0.13	1.35	38,39,41,42	0
5	DMS	A	5020	4/4	0.14	1.30	56,57,57,58	0
5	DMS	D	5017	4/4	0.21	1.28	70,70,71,72	0
5	DMS	C	5015	4/4	0.22	1.25	56,57,57,57	0
5	DMS	D	5028	4/4	0.19	1.25	63,63,63,63	0
5	DMS	A	5126	4/4	0.16	1.23	45,46,46,47	0
5	DMS	D	5016	4/4	0.18	1.09	43,44,44,45	0
5	DMS	B	5029	4/4	0.13	1.08	57,57,58,59	0
5	DMS	D	5012	4/4	0.17	1.07	53,53,53,54	0
3	MG	A	3005	1/1	0.14	1.00	70,70,70,70	0
5	DMS	A	5005	4/4	0.12	0.90	32,32,33,34	0
5	DMS	A	5231	4/4	0.18	0.90	57,57,57,58	0
5	DMS	C	5025	4/4	0.11	0.86	55,55,55,56	0
5	DMS	B	5034	4/4	0.14	0.78	45,46,46,48	0
5	DMS	A	5018	4/4	0.22	0.73	55,55,55,56	0
5	DMS	A	5129	4/4	0.16	0.64	60,61,61,61	0
2	149	B	2001	12/12	0.09	0.59	11,13,16,17	0
5	DMS	A	5003	4/4	0.10	0.49	38,38,38,39	0
5	DMS	C	5005	4/4	0.10	0.43	36,36,37,38	0
5	DMS	B	5008	4/4	0.12	0.43	40,41,41,42	0
5	DMS	A	5008	4/4	0.14	0.38	48,49,49,50	0
5	DMS	A	5232	4/4	0.18	0.35	58,58,58,59	0
5	DMS	B	5024	4/4	0.13	0.22	62,62,62,62	0
5	DMS	B	5012	4/4	0.12	0.18	37,37,38,39	0
5	DMS	A	5012	4/4	0.13	0.17	46,46,47,47	0
5	DMS	B	5005	4/4	0.10	0.16	34,36,37,38	0
5	DMS	D	5005	4/4	0.11	0.15	27,27,28,29	0
5	DMS	D	5025	4/4	0.11	0.13	40,41,41,41	0
2	149	C	2001	12/12	0.08	0.12	9,12,13,15	0
5	DMS	C	5023	4/4	0.10	0.11	47,47,48,48	0
5	DMS	C	5011	4/4	0.12	0.08	41,41,41,42	0
5	DMS	D	5008	4/4	0.11	-0.00	44,45,45,46	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	DMS	A	5023	4/4	0.11	-0.07	60,61,61,62	0
5	DMS	D	5011	4/4	0.11	-0.17	32,32,33,34	0
5	DMS	B	5016	4/4	0.15	-0.19	41,42,42,42	0
5	DMS	A	5130	4/4	0.10	-0.23	27,27,29,31	0
4	NA	C	3102	1/1	0.08	-0.30	14,14,14,14	0
5	DMS	D	5003	4/4	0.09	-0.36	32,33,33,33	0
5	DMS	B	5001	4/4	0.07	-0.36	19,21,23,24	0
5	DMS	C	5010	4/4	0.10	-0.39	34,35,35,36	0
4	NA	C	3104	1/1	0.12	-0.41	32,32,32,32	0
5	DMS	B	5133	4/4	0.12	-0.43	70,70,70,71	0
5	DMS	D	5006	4/4	0.09	-0.54	30,32,33,33	0
4	NA	B	3104	1/1	0.10	-0.72	31,31,31,31	0
4	NA	D	3101	1/1	0.07	-0.74	12,12,12,12	0
4	NA	C	3101	1/1	0.07	-0.77	13,13,13,13	0
5	DMS	A	5001	4/4	0.07	-0.88	22,22,24,24	0
3	MG	D	3001	1/1	0.06	-1.03	13,13,13,13	0
4	NA	A	3101	1/1	0.07	-1.05	13,13,13,13	0
4	NA	B	3103	1/1	0.09	-1.23	23,23,23,23	0
5	DMS	C	5001	4/4	0.07	-1.24	23,25,25,26	0
4	NA	C	3103	1/1	0.09	-1.31	29,29,29,29	0
4	NA	A	3103	1/1	0.08	-1.57	39,39,39,39	0
3	MG	A	3001	1/1	0.07	-1.60	17,17,17,17	0
5	DMS	D	5001	4/4	0.06	-1.63	21,21,22,24	0
4	NA	A	3104	1/1	0.07	-1.67	28,28,28,28	0
4	NA	A	3102	1/1	0.05	-1.82	21,21,21,21	0
4	NA	B	3102	1/1	0.06	-1.98	16,16,16,16	0
4	NA	D	3103	1/1	0.07	-2.05	33,33,33,33	0
3	MG	D	3002	1/1	0.07	-2.14	23,23,23,23	0
3	MG	A	1024	1/1	0.07	-2.22	44,44,44,44	0
4	NA	D	3102	1/1	0.04	-2.84	16,16,16,16	0
5	DMS	A	5011	4/4	0.08	-2.95	39,39,40,41	0
3	MG	C	3002	1/1	0.04	-3.22	17,17,17,17	0
3	MG	B	3001	1/1	0.04	-3.66	15,15,15,15	0
3	MG	A	3002	1/1	0.05	-3.70	19,19,19,19	0
4	NA	B	3101	1/1	0.06	-3.75	15,15,15,15	0
3	MG	B	3002	1/1	0.05	-3.88	19,19,19,19	0
3	MG	C	3001	1/1	0.04	-4.12	12,12,12,12	0

6.5 Other polymers ⓘ

There are no such residues in this entry.