



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 01:14 pm BST

PDB ID : 6MAS
Title : X-ray Structure of Branchiostoma floridae fluorescent protein lanFP10G
Authors : Muslinkina, L.; Pletneva, N.; Pletnev, V.; Pletnev, S.
Deposited on : 2018-08-28
Resolution : 1.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

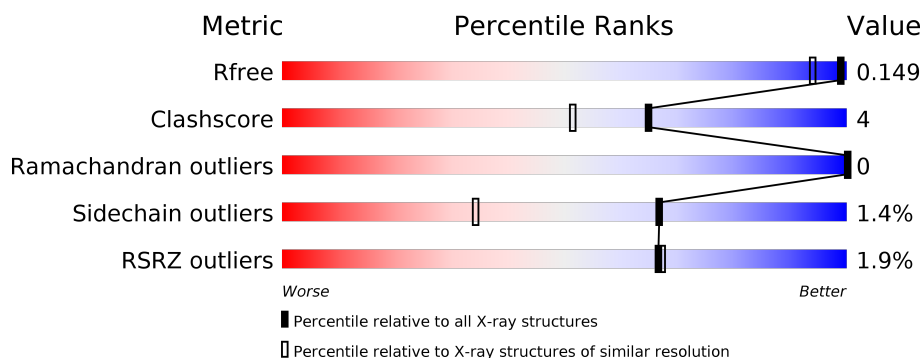
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.30 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}	130704	1058 (1.30-1.30)
Clashscore	141614	1101 (1.30-1.30)
Ramachandran outliers	138981	1058 (1.30-1.30)
Sidechain outliers	138945	1058 (1.30-1.30)
RSRZ outliers	127900	1029 (1.30-1.30)

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 respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	227	<div> <div>3%</div> <div>87%</div> <div>7%</div> <div>5%</div> </div>
1	B	227	<div> <div>%</div> <div>84%</div> <div>10%</div> <div>5%</div> </div>
1	C	227	<div> <div>2%</div> <div>83%</div> <div>11%</div> <div>5%</div> </div>
1	D	227	<div> <div>2%</div> <div>86%</div> <div>8%</div> <div>5%</div> </div>
1	E	227	<div> <div>%</div> <div>90%</div> <div>•</div> <div>5%</div> </div>
1	F	227	<div> <div>2%</div> <div>88%</div> <div>7%</div> <div>5%</div> </div>

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Mol	Chain	Length	Quality of chain
1	G	227	<div><div></div><div>89%</div><div>6%5%</div></div>
1	H	227	<div><div>4%</div><div></div><div>86%</div><div>8%5%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17339 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	16	0
			1878	1199	318	349	12			
1	B	216	Total	C	N	O	S	0	13	0
			1849	1183	308	345	13			
1	C	216	Total	C	N	O	S	0	20	0
			1905	1221	317	352	15			
1	D	216	Total	C	N	O	S	0	16	0
			1874	1198	313	349	14			
1	E	216	Total	C	N	O	S	0	12	0
			1840	1178	303	346	13			
1	F	216	Total	C	N	O	S	0	11	0
			1834	1174	307	340	13			
1	G	216	Total	C	N	O	S	0	9	0
			1813	1159	303	338	13			
1	H	216	Total	C	N	O	S	0	13	0
			1843	1178	305	346	14			

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP C3YRA2
A	58	CR2	GLY	chromophore	UNP C3YRA2
A	58	CR2	TYR	chromophore	UNP C3YRA2
A	58	CR2	ALA	chromophore	UNP C3YRA2
A	220	SER	-	expression tag	UNP C3YRA2
A	221	GLY	-	expression tag	UNP C3YRA2
A	222	GLY	-	expression tag	UNP C3YRA2
A	223	SER	-	expression tag	UNP C3YRA2
A	224	HIS	-	expression tag	UNP C3YRA2
A	225	HIS	-	expression tag	UNP C3YRA2
A	226	HIS	-	expression tag	UNP C3YRA2
A	227	HIS	-	expression tag	UNP C3YRA2
A	228	HIS	-	expression tag	UNP C3YRA2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	229	HIS	-	expression tag	UNP C3YRA2
B	1	MET	-	expression tag	UNP C3YRA2
B	58	CR2	GLY	chromophore	UNP C3YRA2
B	58	CR2	TYR	chromophore	UNP C3YRA2
B	58	CR2	ALA	chromophore	UNP C3YRA2
B	220	SER	-	expression tag	UNP C3YRA2
B	221	GLY	-	expression tag	UNP C3YRA2
B	222	GLY	-	expression tag	UNP C3YRA2
B	223	SER	-	expression tag	UNP C3YRA2
B	224	HIS	-	expression tag	UNP C3YRA2
B	225	HIS	-	expression tag	UNP C3YRA2
B	226	HIS	-	expression tag	UNP C3YRA2
B	227	HIS	-	expression tag	UNP C3YRA2
B	228	HIS	-	expression tag	UNP C3YRA2
B	229	HIS	-	expression tag	UNP C3YRA2
C	1	MET	-	expression tag	UNP C3YRA2
C	58	CR2	GLY	chromophore	UNP C3YRA2
C	58	CR2	TYR	chromophore	UNP C3YRA2
C	58	CR2	ALA	chromophore	UNP C3YRA2
C	220	SER	-	expression tag	UNP C3YRA2
C	221	GLY	-	expression tag	UNP C3YRA2
C	222	GLY	-	expression tag	UNP C3YRA2
C	223	SER	-	expression tag	UNP C3YRA2
C	224	HIS	-	expression tag	UNP C3YRA2
C	225	HIS	-	expression tag	UNP C3YRA2
C	226	HIS	-	expression tag	UNP C3YRA2
C	227	HIS	-	expression tag	UNP C3YRA2
C	228	HIS	-	expression tag	UNP C3YRA2
C	229	HIS	-	expression tag	UNP C3YRA2
D	1	MET	-	expression tag	UNP C3YRA2
D	58	CR2	GLY	chromophore	UNP C3YRA2
D	58	CR2	TYR	chromophore	UNP C3YRA2
D	58	CR2	ALA	chromophore	UNP C3YRA2
D	220	SER	-	expression tag	UNP C3YRA2
D	221	GLY	-	expression tag	UNP C3YRA2
D	222	GLY	-	expression tag	UNP C3YRA2
D	223	SER	-	expression tag	UNP C3YRA2
D	224	HIS	-	expression tag	UNP C3YRA2
D	225	HIS	-	expression tag	UNP C3YRA2
D	226	HIS	-	expression tag	UNP C3YRA2
D	227	HIS	-	expression tag	UNP C3YRA2
D	228	HIS	-	expression tag	UNP C3YRA2

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Chain	Residue	Modelled	Actual	Comment	Reference
D	229	HIS	-	expression tag	UNP C3YRA2
E	1	MET	-	expression tag	UNP C3YRA2
E	58	CR2	GLY	chromophore	UNP C3YRA2
E	58	CR2	TYR	chromophore	UNP C3YRA2
E	58	CR2	ALA	chromophore	UNP C3YRA2
E	220	SER	-	expression tag	UNP C3YRA2
E	221	GLY	-	expression tag	UNP C3YRA2
E	222	GLY	-	expression tag	UNP C3YRA2
E	223	SER	-	expression tag	UNP C3YRA2
E	224	HIS	-	expression tag	UNP C3YRA2
E	225	HIS	-	expression tag	UNP C3YRA2
E	226	HIS	-	expression tag	UNP C3YRA2
E	227	HIS	-	expression tag	UNP C3YRA2
E	228	HIS	-	expression tag	UNP C3YRA2
E	229	HIS	-	expression tag	UNP C3YRA2
F	1	MET	-	expression tag	UNP C3YRA2
F	58	CR2	GLY	chromophore	UNP C3YRA2
F	58	CR2	TYR	chromophore	UNP C3YRA2
F	58	CR2	ALA	chromophore	UNP C3YRA2
F	220	SER	-	expression tag	UNP C3YRA2
F	221	GLY	-	expression tag	UNP C3YRA2
F	222	GLY	-	expression tag	UNP C3YRA2
F	223	SER	-	expression tag	UNP C3YRA2
F	224	HIS	-	expression tag	UNP C3YRA2
F	225	HIS	-	expression tag	UNP C3YRA2
F	226	HIS	-	expression tag	UNP C3YRA2
F	227	HIS	-	expression tag	UNP C3YRA2
F	228	HIS	-	expression tag	UNP C3YRA2
F	229	HIS	-	expression tag	UNP C3YRA2
G	1	MET	-	expression tag	UNP C3YRA2
G	58	CR2	GLY	chromophore	UNP C3YRA2
G	58	CR2	TYR	chromophore	UNP C3YRA2
G	58	CR2	ALA	chromophore	UNP C3YRA2
G	220	SER	-	expression tag	UNP C3YRA2
G	221	GLY	-	expression tag	UNP C3YRA2
G	222	GLY	-	expression tag	UNP C3YRA2
G	223	SER	-	expression tag	UNP C3YRA2
G	224	HIS	-	expression tag	UNP C3YRA2
G	225	HIS	-	expression tag	UNP C3YRA2
G	226	HIS	-	expression tag	UNP C3YRA2
G	227	HIS	-	expression tag	UNP C3YRA2
G	228	HIS	-	expression tag	UNP C3YRA2

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Chain	Residue	Modelled	Actual	Comment	Reference
G	229	HIS	-	expression tag	UNP C3YRA2
H	1	MET	-	expression tag	UNP C3YRA2
H	58	CR2	GLY	chromophore	UNP C3YRA2
H	58	CR2	TYR	chromophore	UNP C3YRA2
H	58	CR2	ALA	chromophore	UNP C3YRA2
H	220	SER	-	expression tag	UNP C3YRA2
H	221	GLY	-	expression tag	UNP C3YRA2
H	222	GLY	-	expression tag	UNP C3YRA2
H	223	SER	-	expression tag	UNP C3YRA2
H	224	HIS	-	expression tag	UNP C3YRA2
H	225	HIS	-	expression tag	UNP C3YRA2
H	226	HIS	-	expression tag	UNP C3YRA2
H	227	HIS	-	expression tag	UNP C3YRA2
H	228	HIS	-	expression tag	UNP C3YRA2
H	229	HIS	-	expression tag	UNP C3YRA2

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	E	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	F	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		
2	G	1	Total	C	O	0	0
			6	3	3		

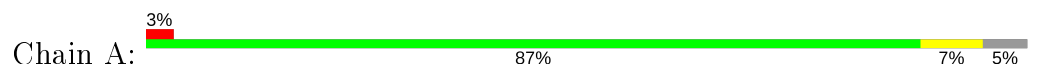
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	311	Total	O	0	0
			311	311		
3	B	321	Total	O	0	0
			321	321		
3	C	331	Total	O	0	0
			331	331		
3	D	322	Total	O	0	0
			322	322		
3	E	295	Total	O	0	0
			295	295		
3	F	300	Total	O	0	0
			300	300		
3	G	299	Total	O	0	0
			299	299		
3	H	282	Total	O	0	0
			282	282		

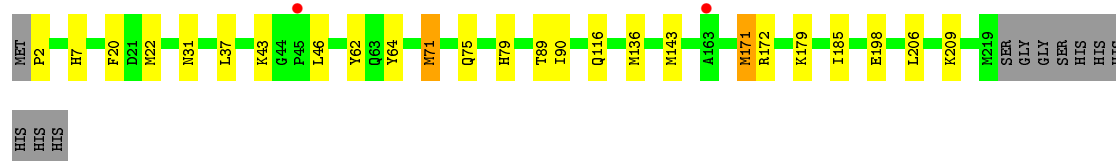
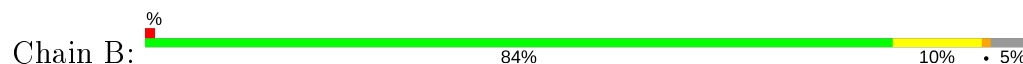
3 Residue-property plots [i](#)

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

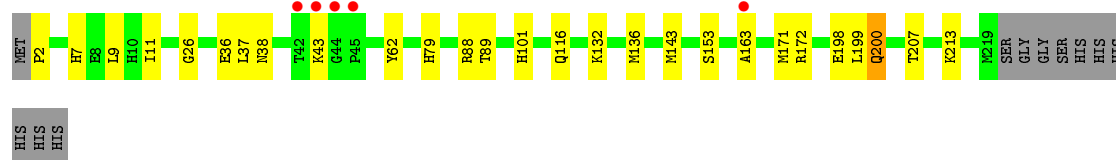
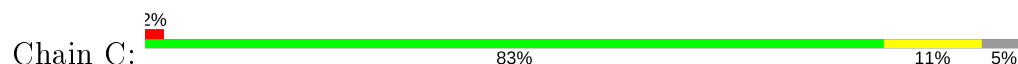
- Molecule 1: Uncharacterized protein



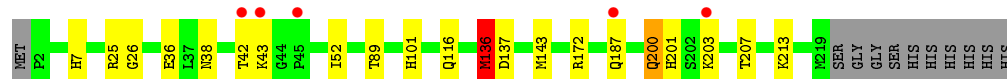
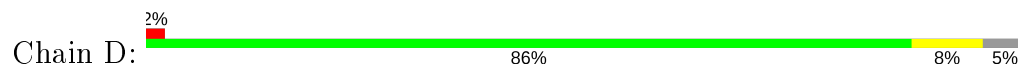
- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein

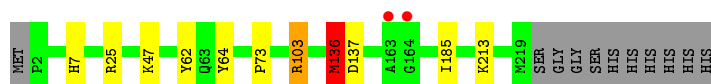


- Molecule 1: Uncharacterized protein

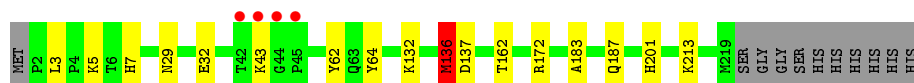
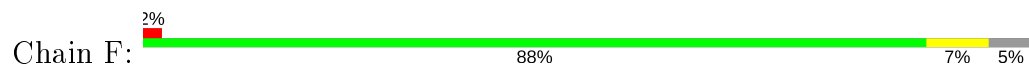


- Molecule 1: Uncharacterized protein

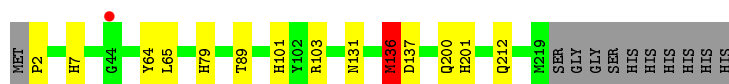




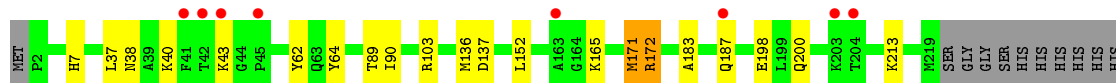
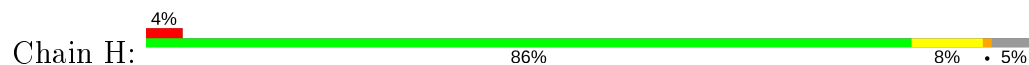
- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	77.88Å 104.31Å 116.14Å 90.00° 95.32° 90.00°	Depositor
Resolution (Å)	29.82 – 1.30 29.80 – 1.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.82-1.30) 99.9 (29.80-1.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 1.30Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.119 , 0.147 0.122 , 0.149	Depositor DCC
R_{free} test set	4625 reflections (1.03%)	wwPDB-VP
Wilson B-factor (Å ²)	11.3	Xtriage
Anisotropy	0.235	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 49.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	17339	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CR2, GOL

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	0.55	0/1910	0.80	1/2573 (0.0%)
1	B	0.59	0/1882	0.91	9/2540 (0.4%)
1	C	0.58	0/1938	0.87	6/2613 (0.2%)
1	D	0.56	0/1907	0.83	3/2574 (0.1%)
1	E	0.56	0/1873	0.83	7/2529 (0.3%)
1	F	0.57	0/1867	0.84	6/2519 (0.2%)
1	G	0.56	0/1845	0.78	2/2490 (0.1%)
1	H	0.57	0/1875	0.85	5/2532 (0.2%)
All	All	0.57	0/15097	0.84	39/20370 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1
1	E	0	1
1	G	0	1
1	H	0	1
All	All	0	4

There are no bond length outliers.

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	136[A]	MET	CG-SD-CE	-10.77	82.96	100.20
1	C	136[B]	MET	CG-SD-CE	-10.77	82.96	100.20
1	B	62	TYR	CB-CG-CD2	-8.27	116.04	121.00
1	B	62	TYR	CB-CG-CD1	6.80	125.08	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	103	ARG	CG-CD-NE	-6.62	97.89	111.80
1	B	143	MET	CG-SD-CE	-6.62	89.61	100.20
1	H	103	ARG	CG-CD-NE	-6.61	97.93	111.80
1	B	171[A]	MET	CG-SD-CE	-6.11	90.43	100.20
1	B	171[B]	MET	CG-SD-CE	-6.11	90.43	100.20
1	A	143	MET	CG-SD-CE	-6.10	90.44	100.20
1	H	62	TYR	CB-CG-CD2	-6.10	117.34	121.00
1	C	62[A]	TYR	CB-CG-CD2	-6.03	117.38	121.00
1	C	62[B]	TYR	CB-CG-CD2	-6.03	117.38	121.00
1	B	136	MET	CG-SD-CE	-6.03	90.55	100.20
1	F	62[A]	TYR	CB-CG-CD2	-5.96	117.42	121.00
1	F	62[B]	TYR	CB-CG-CD2	-5.96	117.42	121.00
1	B	71[A]	MET	CG-SD-CE	-5.94	90.69	100.20
1	B	71[B]	MET	CG-SD-CE	-5.94	90.69	100.20
1	G	136[A]	MET	CG-SD-CE	-5.94	90.69	100.20
1	G	136[B]	MET	CG-SD-CE	-5.94	90.69	100.20
1	H	171[A]	MET	CG-SD-CE	-5.90	90.76	100.20
1	H	171[B]	MET	CG-SD-CE	-5.90	90.76	100.20
1	D	136[A]	MET	CG-SD-CE	-5.90	90.76	100.20
1	D	136[B]	MET	CG-SD-CE	-5.90	90.76	100.20
1	E	62[A]	TYR	CB-CG-CD2	-5.48	117.72	121.00
1	E	62[B]	TYR	CB-CG-CD2	-5.48	117.72	121.00
1	F	62[A]	TYR	CB-CG-CD1	5.47	124.28	121.00
1	F	62[B]	TYR	CB-CG-CD1	5.47	124.28	121.00
1	D	143	MET	CG-SD-CE	-5.45	91.48	100.20
1	E	136[A]	MET	CG-SD-CE	-5.38	91.59	100.20
1	E	136[B]	MET	CG-SD-CE	-5.38	91.59	100.20
1	B	22	MET	CG-SD-CE	-5.35	91.64	100.20
1	C	62[A]	TYR	CB-CG-CD1	5.35	124.21	121.00
1	C	62[B]	TYR	CB-CG-CD1	5.35	124.21	121.00
1	F	136[A]	MET	CG-SD-CE	-5.26	91.78	100.20
1	F	136[B]	MET	CG-SD-CE	-5.26	91.78	100.20
1	H	103	ARG	NE-CZ-NH1	-5.09	117.76	120.30
1	E	62[A]	TYR	CB-CG-CD1	5.03	124.02	121.00
1	E	62[B]	TYR	CB-CG-CD1	5.03	124.02	121.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	172[A]	ARG	Sidechain
1	E	103	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	G	103[A]	ARG	Sidechain
1	H	172[A]	ARG	Sidechain

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1878	0	1810	14	0
1	B	1849	0	1766	20	0
1	C	1905	0	1838	28	0
1	D	1874	0	1796	17	0
1	E	1840	0	1751	7	0
1	F	1834	0	1761	11	0
1	G	1813	0	1736	11	0
1	H	1843	0	1760	13	0
2	B	6	0	8	0	0
2	C	6	0	8	0	0
2	D	6	0	8	0	0
2	E	6	0	8	0	0
2	F	12	0	16	0	0
2	G	6	0	8	0	0
3	A	311	0	0	7	0
3	B	321	0	0	7	0
3	C	331	0	0	8	0
3	D	322	0	0	11	0
3	E	295	0	0	3	0
3	F	300	0	0	6	0
3	G	299	0	0	5	0
3	H	282	0	0	6	0
All	All	17339	0	14274	119	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (119) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:88:ARG:HD3	1:C:171[A]:MET:CE	1.81	1.09
1:C:88:ARG:HD3	1:C:171[A]:MET:HE3	1.29	1.06
1:E:47:LYS:HE2	3:E:628:HOH:O	1.61	0.99
1:C:88:ARG:NE	1:C:171[A]:MET:HE2	1.79	0.98
1:B:172[A]:ARG:NH1	3:B:401:HOH:O	1.94	0.97
1:C:88:ARG:CD	1:C:171[A]:MET:CE	2.45	0.93
1:A:116:GLN:HE22	1:B:116:GLN:HE22	1.20	0.90
1:C:88:ARG:CD	1:C:171[A]:MET:HE3	2.02	0.89
1:C:88:ARG:NE	1:C:171[A]:MET:CE	2.38	0.86
1:C:88:ARG:CZ	1:C:171[A]:MET:CE	2.55	0.84
1:C:88:ARG:CD	1:C:171[A]:MET:HE2	2.08	0.83
1:F:172[A]:ARG:NH2	3:F:401:HOH:O	1.95	0.83
1:C:116:GLN:HE22	1:D:116:GLN:HE22	1.29	0.80
1:B:20[B]:PHE:HE1	1:B:46:LEU:HD21	1.46	0.79
1:C:38[B]:ASN:ND2	1:C:207[B]:THR:HG22	2.02	0.74
1:A:22:MET:HE1	1:A:53:LEU:HD13	1.70	0.72
1:D:207[B]:THR:HG23	3:D:493:HOH:O	1.89	0.72
1:A:40[B]:LYS:HG2	3:A:348:HOH:O	1.93	0.69
1:B:89[B]:THR:HG23	3:B:564:HOH:O	1.93	0.69
1:C:89[A]:THR:HG21	3:D:401:HOH:O	1.92	0.68
1:B:198[B]:GLU:HG2	1:B:209:LYS:HB2	1.75	0.66
1:B:20[B]:PHE:HE1	1:B:46:LEU:CD2	2.09	0.66
1:D:101:HIS:HD2	3:D:604:HOH:O	1.77	0.66
1:D:89[B]:THR:HG23	3:D:591:HOH:O	1.97	0.65
1:A:47:LYS:NZ	3:A:301:HOH:O	2.31	0.63
1:A:89:THR:HG21	3:B:405:HOH:O	1.97	0.62
1:E:213[B]:LYS:HE3	3:E:658:HOH:O	2.00	0.60
1:H:183:ALA:O	1:H:187:GLN:HG3	2.01	0.60
3:A:429:HOH:O	1:B:89[A]:THR:HG21	2.01	0.60
1:B:2:PRO:N	1:B:79:HIS:HD1	2.01	0.59
1:F:213[B]:LYS:HE2	3:F:552:HOH:O	2.02	0.59
1:D:38:ASN:HD22	1:D:207[B]:THR:HG22	1.67	0.59
1:G:89:THR:HG21	3:H:307:HOH:O	2.02	0.58
1:F:132:LYS:HB3	1:F:162[B]:THR:HG22	1.86	0.58
1:G:2:PRO:N	1:G:79:HIS:HD1	2.01	0.58
1:H:89[B]:THR:HG23	3:H:459:HOH:O	2.05	0.57
1:H:7:HIS:HE1	3:H:336:HOH:O	1.89	0.56
1:A:22:MET:HE1	1:A:53:LEU:CD1	2.35	0.55
1:C:163:ALA:N	3:C:407:HOH:O	2.38	0.55
1:D:136[B]:MET:HG3	1:D:137:ASP:N	2.22	0.55
1:D:7:HIS:HE1	3:D:429:HOH:O	1.90	0.54
1:G:201:HIS:HD2	3:G:405:HOH:O	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:90:ILE:HG12	1:H:171[A]:MET:HG3	1.90	0.54
1:C:88:ARG:CZ	1:C:171[A]:MET:HE1	2.38	0.54
1:C:2:PRO:N	1:C:79:HIS:HD1	2.06	0.53
1:B:7:HIS:HE1	3:B:423:HOH:O	1.90	0.53
1:H:213:LYS:HE3	3:H:537:HOH:O	2.07	0.53
1:C:38[B]:ASN:HD22	1:C:207[B]:THR:HG22	1.74	0.53
1:C:9[B]:LEU:HD21	1:C:11[B]:ILE:HG13	1.91	0.52
1:C:7:HIS:HE1	3:C:444:HOH:O	1.93	0.52
1:C:132:LYS:HD3	3:C:407:HOH:O	2.08	0.51
1:G:7:HIS:HE1	3:G:426:HOH:O	1.92	0.51
1:F:201:HIS:HE1	3:F:621:HOH:O	1.94	0.51
1:H:136[A]:MET:HG3	1:H:137:ASP:N	2.25	0.51
1:B:20[B]:PHE:CE1	1:B:46:LEU:HD21	2.36	0.50
1:A:7:HIS:HE1	3:A:331:HOH:O	1.94	0.50
1:F:5:LYS:HD2	3:F:572:HOH:O	2.11	0.50
1:A:40[A]:LYS:HG2	1:A:205:GLU:HB3	1.94	0.50
1:F:7:HIS:HE1	3:F:440:HOH:O	1.94	0.49
1:A:156:ASP:OD1	1:A:170[B]:LYS:HD2	2.12	0.49
1:E:7:HIS:HE1	3:E:443:HOH:O	1.95	0.49
1:A:2:PRO:N	1:A:79:HIS:HD1	2.10	0.49
1:C:213[B]:LYS:HE2	3:C:414:HOH:O	2.12	0.49
1:A:136:MET:CE	3:A:314:HOH:O	2.61	0.48
1:B:71[A]:MET:HG2	1:B:75:GLN:HB3	1.94	0.48
1:F:183:ALA:O	1:F:187:GLN:HG3	2.14	0.48
1:C:199[A]:LEU:HD22	1:C:207[A]:THR:O	2.14	0.47
1:A:22:MET:HE2	1:A:22:MET:HB2	1.38	0.47
1:B:71[A]:MET:HG2	1:B:75:GLN:CB	2.45	0.47
1:F:136[B]:MET:HG3	1:F:137:ASP:N	2.30	0.47
1:H:152:LEU:HD21	1:H:172[B]:ARG:NH2	2.30	0.47
1:D:201:HIS:HE1	3:D:632:HOH:O	1.98	0.47
1:E:136[B]:MET:HG3	1:E:137:ASP:N	2.30	0.46
1:B:179:LYS:NZ	3:B:403:HOH:O	2.24	0.46
1:D:187:GLN:HG3	3:D:546:HOH:O	2.16	0.46
1:C:200:GLN:NE2	3:C:410:HOH:O	2.50	0.45
1:D:207[A]:THR:HG23	3:D:539:HOH:O	2.15	0.45
1:B:31:ASN:ND2	3:B:406:HOH:O	2.44	0.45
1:D:213[B]:LYS:HE3	3:D:564:HOH:O	2.17	0.45
1:C:143[A]:MET:HG2	1:C:153:SER:OG	2.16	0.45
1:D:42:THR:HG22	1:D:43:LYS:HG3	1.99	0.44
1:G:131:ASN:ND2	3:G:403:HOH:O	2.40	0.44
1:G:200[A]:GLN:HA	1:G:200[A]:GLN:OE1	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:ILE:HG12	1:B:171[A]:MET:HG3	1.99	0.44
1:B:89[B]:THR:HG22	3:B:543:HOH:O	2.18	0.43
1:D:200:GLN:NE2	3:D:402:HOH:O	2.26	0.43
3:A:484:HOH:O	1:C:172[A]:ARG:HD2	2.18	0.43
1:G:101:HIS:HE1	3:G:502:HOH:O	2.01	0.43
1:H:172[B]:ARG:NH1	3:H:305:HOH:O	2.48	0.43
1:A:149[A]:LYS:HE3	3:A:328:HOH:O	2.18	0.43
1:H:198[B]:GLU:CG	1:H:200:GLN:HE22	2.32	0.43
1:B:7:HIS:HD2	1:B:64:TYR:OH	2.02	0.43
1:D:203:LYS:HA	1:D:203:LYS:HD3	1.88	0.43
1:F:3:LEU:HD23	3:F:536:HOH:O	2.18	0.43
1:C:198[B]:GLU:HG2	1:C:200:GLN:OE1	2.19	0.43
1:C:101:HIS:HE1	3:C:530:HOH:O	2.02	0.43
1:B:7:HIS:CD2	1:B:64:TYR:OH	2.72	0.42
1:D:207[A]:THR:CG2	3:D:539:HOH:O	2.67	0.42
1:G:136[B]:MET:HG3	1:G:137:ASP:N	2.33	0.42
1:H:198[B]:GLU:CD	1:H:200:GLN:HE22	2.22	0.42
1:D:26:GLY:HA3	1:D:36:GLU:O	2.19	0.42
1:G:136[B]:MET:HE1	3:G:682:HOH:O	2.20	0.41
1:C:200:GLN:HB2	1:C:200:GLN:HE21	1.67	0.41
1:G:7:HIS:CD2	1:G:64:TYR:OH	2.74	0.41
1:D:52:ILE:HD12	1:D:52:ILE:HA	1.95	0.41
1:C:200:GLN:NE2	3:C:403:HOH:O	2.33	0.41
1:G:65:LEU:HD13	1:G:212:GLN:HB2	2.02	0.41
1:E:7:HIS:HD2	1:E:64:TYR:OH	2.03	0.41
1:A:172[A]:ARG:HD2	3:C:590:HOH:O	2.21	0.41
1:B:71[B]:MET:HE2	1:B:75:GLN:HB3	2.03	0.41
1:E:73:PRO:HD3	1:E:185[A]:ILE:HD12	2.03	0.41
1:E:7:HIS:CD2	1:E:64:TYR:OH	2.74	0.41
1:F:29:ASN:ND2	1:F:32:GLU:HG3	2.37	0.41
1:H:38:ASN:HB2	3:H:517:HOH:O	2.20	0.40
1:B:206:LEU:C	1:B:206:LEU:HD12	2.41	0.40
1:C:26:GLY:HA3	1:C:36:GLU:O	2.21	0.40
1:H:198[B]:GLU:HG2	1:H:200:GLN:NE2	2.36	0.40
1:H:7:HIS:HD2	1:H:64:TYR:OH	2.05	0.40
1:F:7:HIS:CD2	1:F:64:TYR:OH	2.74	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	227/227 (100%)	226 (100%)	1 (0%)	0	100	100
1	B	224/227 (99%)	223 (100%)	1 (0%)	0	100	100
1	C	231/227 (102%)	229 (99%)	2 (1%)	0	100	100
1	D	227/227 (100%)	225 (99%)	2 (1%)	0	100	100
1	E	223/227 (98%)	221 (99%)	2 (1%)	0	100	100
1	F	222/227 (98%)	220 (99%)	2 (1%)	0	100	100
1	G	220/227 (97%)	220 (100%)	0	0	100	100
1	H	224/227 (99%)	222 (99%)	2 (1%)	0	100	100
All	All	1798/1816 (99%)	1786 (99%)	12 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	202/195 (104%)	199 (98%)	3 (2%)	65	31
1	B	199/195 (102%)	196 (98%)	3 (2%)	65	31
1	C	206/195 (106%)	203 (98%)	3 (2%)	65	31
1	D	202/195 (104%)	197 (98%)	5 (2%)	47	10
1	E	198/195 (102%)	195 (98%)	3 (2%)	65	31
1	F	197/195 (101%)	194 (98%)	3 (2%)	65	31

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	195/195 (100%)	193 (99%)	2 (1%)	76	48
1	H	199/195 (102%)	195 (98%)	4 (2%)	55	17
All	All	1598/1560 (102%)	1572 (98%)	26 (2%)	67	28

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	LEU
1	A	188	LYS
1	A	205	GLU
1	B	37	LEU
1	B	43	LYS
1	B	185	ILE
1	C	37	LEU
1	C	43	LYS
1	C	200	GLN
1	D	25[A]	ARG
1	D	25[B]	ARG
1	D	136[A]	MET
1	D	136[B]	MET
1	D	200	GLN
1	E	25	ARG
1	E	136[A]	MET
1	E	136[B]	MET
1	F	43	LYS
1	F	136[A]	MET
1	F	136[B]	MET
1	G	136[A]	MET
1	G	136[B]	MET
1	H	37	LEU
1	H	40	LYS
1	H	43	LYS
1	H	165	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	HIS
1	A	101	HIS
1	A	131	ASN

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Mol	Chain	Res	Type
1	A	200	GLN
1	A	201	HIS
1	B	7	HIS
1	B	101	HIS
1	B	116	GLN
1	C	7	HIS
1	C	101	HIS
1	C	168	GLN
1	C	175	ASN
1	C	201	HIS
1	D	7	HIS
1	D	38	ASN
1	D	101	HIS
1	D	116	GLN
1	D	201	HIS
1	E	7	HIS
1	E	79	HIS
1	E	85	GLN
1	E	101	HIS
1	E	116	GLN
1	F	7	HIS
1	F	101	HIS
1	F	201	HIS
1	G	7	HIS
1	G	101	HIS
1	G	116	GLN
1	G	168	GLN
1	G	201	HIS
1	H	7	HIS
1	H	101	HIS
1	H	168	GLN
1	H	201	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
1	CR2	B	58	1	20,20,21	2.50	3 (15%)	25,27,29	1.50	6 (24%)
1	CR2	C	58	1	20,20,21	2.22	3 (15%)	25,27,29	1.71	6 (24%)
1	CR2	A	58	1	20,20,21	2.78	4 (20%)	25,27,29	2.12	6 (24%)
1	CR2	F	58	1	20,20,21	3.42	4 (20%)	25,27,29	2.99	9 (36%)
1	CR2	G	58	1	20,20,21	2.67	4 (20%)	25,27,29	2.02	4 (16%)
1	CR2	D	58	1	20,20,21	3.11	5 (25%)	25,27,29	3.06	9 (36%)
1	CR2	E	58	1	20,20,21	3.04	4 (20%)	25,27,29	2.78	5 (20%)
1	CR2	H	58	1	20,20,21	2.88	3 (15%)	25,27,29	2.85	9 (36%)

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
1	CR2	B	58	1	-	0/6/25/26	0/2/2/2
1	CR2	C	58	1	-	0/6/25/26	0/2/2/2
1	CR2	A	58	1	-	0/6/25/26	0/2/2/2
1	CR2	F	58	1	-	0/6/25/26	0/2/2/2
1	CR2	G	58	1	-	0/6/25/26	0/2/2/2
1	CR2	D	58	1	-	0/6/25/26	0/2/2/2
1	CR2	E	58	1	-	0/6/25/26	0/2/2/2
1	CR2	H	58	1	-	0/6/25/26	0/2/2/2

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	58	CR2	CB2-CA2	11.74	1.44	1.35
1	D	58	CR2	CB2-CA2	10.47	1.43	1.35
1	E	58	CR2	CB2-CA2	10.16	1.43	1.35
1	A	58	CR2	CB2-CA2	10.15	1.43	1.35
1	H	58	CR2	CB2-CA2	9.37	1.42	1.35
1	B	58	CR2	CB2-CA2	9.28	1.42	1.35
1	G	58	CR2	CB2-CA2	8.87	1.42	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	58	CR2	CA2-C2	-8.27	1.40	1.48
1	C	58	CR2	CB2-CA2	7.42	1.41	1.35
1	H	58	CR2	CA2-C2	-7.39	1.41	1.48
1	E	58	CR2	CA2-C2	-7.29	1.41	1.48
1	D	58	CR2	CA2-C2	-7.02	1.41	1.48
1	G	58	CR2	CA2-C2	-6.33	1.42	1.48
1	A	58	CR2	CA2-C2	-5.40	1.43	1.48
1	C	58	CR2	CA2-C2	-4.97	1.43	1.48
1	B	58	CR2	CA2-C2	-4.02	1.44	1.48
1	E	58	CR2	C2-N3	-3.53	1.31	1.39
1	B	58	CR2	OH-CZ	-3.17	1.29	1.37
1	F	58	CR2	OH-CZ	-3.16	1.29	1.37
1	D	58	CR2	O2-C2	2.91	1.29	1.23
1	D	58	CR2	OH-CZ	-2.81	1.30	1.37
1	A	58	CR2	OH-CZ	-2.77	1.30	1.37
1	F	58	CR2	C2-N3	-2.68	1.33	1.39
1	G	58	CR2	C2-N3	-2.65	1.33	1.39
1	C	58	CR2	C2-N3	-2.60	1.33	1.39
1	E	58	CR2	OH-CZ	-2.56	1.31	1.37
1	D	58	CR2	C2-N3	-2.44	1.34	1.39
1	H	58	CR2	CG2-CB2	-2.21	1.42	1.46
1	G	58	CR2	OH-CZ	-2.12	1.32	1.37
1	A	58	CR2	C2-N3	-2.06	1.35	1.39

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	58	CR2	CA2-C2-N3	11.19	108.66	103.37
1	F	58	CR2	CA2-C2-N3	10.44	108.31	103.37
1	H	58	CR2	CA2-C2-N3	9.38	107.81	103.37
1	E	58	CR2	CA2-C2-N3	9.16	107.70	103.37
1	E	58	CR2	O2-C2-CA2	-8.81	126.01	130.96
1	H	58	CR2	O2-C2-CA2	-7.35	126.83	130.96
1	F	58	CR2	O2-C2-CA2	-7.11	126.97	130.96
1	A	58	CR2	CA2-C2-N3	7.03	106.69	103.37
1	G	58	CR2	CA2-C2-N3	6.62	106.50	103.37
1	D	58	CR2	O2-C2-CA2	-5.97	127.61	130.96
1	G	58	CR2	O2-C2-CA2	-5.03	128.13	130.96
1	C	58	CR2	CA2-C2-N3	4.50	105.50	103.37
1	A	58	CR2	O2-C2-CA2	-4.27	128.56	130.96
1	D	58	CR2	C2-CA2-N2	-3.77	106.29	108.93
1	A	58	CR2	O3-C3-CA3	-3.76	115.03	126.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	58	CR2	O2-C2-CA2	-3.75	128.85	130.96
1	F	58	CR2	O3-C3-CA3	-3.68	115.29	126.39
1	H	58	CR2	O3-C3-CA3	-3.54	115.69	126.39
1	G	58	CR2	O3-C3-CA3	-3.39	116.14	126.39
1	B	58	CR2	O3-C3-CA3	-3.36	116.24	126.39
1	C	58	CR2	O3-C3-CA3	-3.34	116.31	126.39
1	D	58	CR2	O3-C3-CA3	-3.24	116.62	126.39
1	B	58	CR2	CA2-C2-N3	3.23	104.90	103.37
1	H	58	CR2	C1-CA1-N1	-3.17	105.83	112.85
1	D	58	CR2	C1-CA1-N1	-3.07	106.05	112.85
1	F	58	CR2	C1-CA1-N1	-2.96	106.31	112.85
1	B	58	CR2	C1-CA1-N1	-2.89	106.47	112.85
1	D	58	CR2	CA1-C1-N3	2.84	126.33	122.52
1	H	58	CR2	C2-CA2-N2	-2.68	107.06	108.93
1	G	58	CR2	C1-CA1-N1	-2.67	106.94	112.85
1	A	58	CR2	C1-CA1-N1	-2.62	107.05	112.85
1	F	58	CR2	CA1-C1-N3	2.50	125.87	122.52
1	E	58	CR2	C1-CA1-N1	-2.50	107.32	112.85
1	B	58	CR2	O2-C2-CA2	-2.49	129.56	130.96
1	F	58	CR2	C2-N3-C1	-2.47	106.78	107.99
1	E	58	CR2	O3-C3-CA3	-2.42	119.08	126.39
1	A	58	CR2	C2-CA2-N2	-2.42	107.24	108.93
1	D	58	CR2	CD2-CE2-CZ	2.41	122.52	119.88
1	C	58	CR2	CG2-CB2-CA2	2.39	132.87	129.94
1	C	58	CR2	CA1-C1-N3	2.36	125.68	122.52
1	C	58	CR2	C1-CA1-N1	-2.36	107.64	112.85
1	F	58	CR2	CD2-CE2-CZ	2.35	122.45	119.88
1	H	58	CR2	CA1-C1-N3	2.31	125.61	122.52
1	B	58	CR2	CA1-C1-N3	2.19	125.45	122.52
1	D	58	CR2	CD1-CE1-CZ	2.15	122.24	119.88
1	H	58	CR2	CD1-CE1-CZ	2.15	122.23	119.88
1	F	58	CR2	C2-CA2-N2	-2.10	107.46	108.93
1	B	58	CR2	CD1-CE1-CZ	2.08	122.16	119.88
1	D	58	CR2	CE2-CZ-CE1	-2.07	116.28	119.77
1	A	58	CR2	CA1-C1-N3	2.06	125.28	122.52
1	H	58	CR2	CE2-CZ-CE1	-2.03	116.34	119.77
1	H	58	CR2	CD2-CE2-CZ	2.03	122.10	119.88
1	E	58	CR2	CA1-C1-N3	2.01	125.21	122.52
1	F	58	CR2	CD1-CE1-CZ	2.01	122.08	119.88

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

7 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	GOL	F	302	-	5,5,5	0.64	0	5,5,5	1.54	0
2	GOL	F	301	-	5,5,5	0.32	0	5,5,5	2.33	2 (40%)
2	GOL	D	301	-	5,5,5	0.54	0	5,5,5	0.69	0
2	GOL	G	301	-	5,5,5	1.87	1 (20%)	5,5,5	2.12	2 (40%)
2	GOL	B	301	-	5,5,5	1.47	1 (20%)	5,5,5	1.75	1 (20%)
2	GOL	E	301	-	5,5,5	0.68	0	5,5,5	0.41	0
2	GOL	C	301	-	5,5,5	1.17	0	5,5,5	1.32	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	GOL	F	302	-	-	2/4/4/4	-
2	GOL	F	301	-	-	2/4/4/4	-
2	GOL	D	301	-	-	0/4/4/4	-
2	GOL	G	301	-	-	2/4/4/4	-
2	GOL	B	301	-	-	2/4/4/4	-
2	GOL	E	301	-	-	0/4/4/4	-
2	GOL	C	301	-	-	0/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	301	GOL	O2-C2	3.52	1.53	1.43
2	B	301	GOL	O2-C2	2.91	1.52	1.43

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	301	GOL	O3-C3-C2	4.34	131.03	110.20
2	B	301	GOL	C3-C2-C1	-3.35	98.67	111.70
2	G	301	GOL	O2-C2-C1	3.02	122.41	109.12
2	G	301	GOL	C3-C2-C1	-2.88	100.50	111.70
2	F	301	GOL	O2-C2-C1	-2.28	99.07	109.12

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	302	GOL	O1-C1-C2-C3
2	F	302	GOL	O1-C1-C2-O2
2	F	301	GOL	O1-C1-C2-C3
2	G	301	GOL	O1-C1-C2-C3
2	B	301	GOL	O1-C1-C2-C3
2	B	301	GOL	O1-C1-C2-O2
2	F	301	GOL	O1-C1-C2-O2
2	G	301	GOL	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2			OWAB(Å ²)	Q < 0.9
1	A	215/227 (94%)	-0.40	6 (2%)	53	50	7, 12, 28, 45	0
1	B	215/227 (94%)	-0.51	2 (0%)	84	85	8, 12, 27, 40	0
1	C	215/227 (94%)	-0.42	5 (2%)	60	59	6, 10, 26, 47	0
1	D	215/227 (94%)	-0.42	5 (2%)	60	59	7, 11, 25, 44	0
1	E	215/227 (94%)	-0.50	2 (0%)	84	85	9, 13, 26, 38	0
1	F	215/227 (94%)	-0.45	4 (1%)	66	67	9, 13, 29, 54	0
1	G	215/227 (94%)	-0.58	1 (0%)	91	91	8, 13, 27, 42	0
1	H	215/227 (94%)	-0.24	8 (3%)	41	38	9, 14, 31, 52	0
All	All	1720/1816 (94%)	-0.44	33 (1%)	66	67	6, 13, 28, 54	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	42	THR	6.1
1	D	42	THR	5.8
1	A	204	THR	4.6
1	F	42	THR	4.5
1	F	45	PRO	4.1
1	H	43	LYS	4.1
1	H	41	PHE	4.0
1	C	42	THR	3.7
1	B	163	ALA	3.6
1	C	44	GLY	3.6
1	A	203	LYS	3.4
1	C	43	LYS	3.3
1	H	203	LYS	3.3
1	A	163	ALA	3.2
1	H	187	GLN	3.1
1	A	45	PRO	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	203	LYS	2.9
1	F	43	LYS	2.9
1	A	42	THR	2.8
1	D	43	LYS	2.8
1	E	163	ALA	2.8
1	H	163	ALA	2.7
1	G	44	GLY	2.6
1	D	45	PRO	2.6
1	B	45	PRO	2.5
1	C	163	ALA	2.5
1	E	164	GLY	2.5
1	D	187	GLN	2.4
1	H	204	THR	2.4
1	H	45	PRO	2.3
1	F	44	GLY	2.1
1	C	45	PRO	2.1
1	A	47	LYS	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	CR2	B	58	19/20	0.98	0.07	8,9,11,11	0
1	CR2	G	58	19/20	0.98	0.06	8,10,12,12	0
1	CR2	H	58	19/20	0.98	0.08	10,11,12,12	0
1	CR2	F	58	19/20	0.99	0.05	10,10,11,11	0
1	CR2	C	58	19/20	0.99	0.07	6,7,9,9	0
1	CR2	D	58	19/20	0.99	0.05	7,8,10,10	0
1	CR2	E	58	19/20	0.99	0.07	9,10,11,12	0
1	CR2	A	58	19/20	0.99	0.07	7,8,9,9	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	GOL	F	302	6/6	0.53	0.31	24,26,31,32	0
2	GOL	F	301	6/6	0.75	0.20	23,24,29,30	0
2	GOL	C	301	6/6	0.83	0.23	20,21,22,24	0
2	GOL	B	301	6/6	0.87	0.16	20,23,26,28	0
2	GOL	D	301	6/6	0.90	0.21	20,23,26,26	0
2	GOL	G	301	6/6	0.90	0.12	23,25,27,27	0
2	GOL	E	301	6/6	0.98	0.05	15,15,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.