



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 4, 2021 – 02:19 PM EDT

PDB ID : 7RRI
Title : Crystal structure of fast switching S142A/M159T mutant of fluorescent protein Dronpa (Dronpa2)
Authors : Lin, C.-Y.; Romei, M.G.; Mathews, I.I.; Boxer, S.G.
Deposited on : 2021-08-09
Resolution : 2.64 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 : 2.23.2
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) : 2.23.2

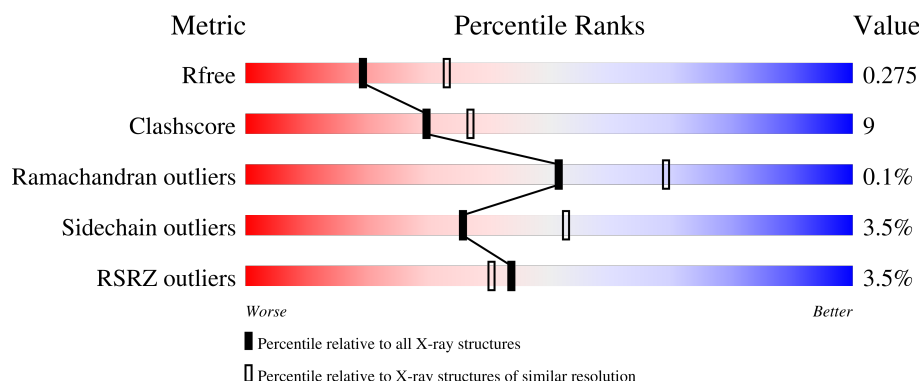
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.64 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	1426 (2.66-2.62)
Clashscore	141614	1472 (2.66-2.62)
Ramachandran outliers	138981	1446 (2.66-2.62)
Sidechain outliers	138945	1446 (2.66-2.62)
RSRZ outliers	127900	1408 (2.66-2.62)

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 of 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	255	<div> <div>2%</div> <div> <div></div> <div>67%</div> <div>16%</div> <div>•</div> <div>15%</div> </div> </div>
1	B	255	<div> <div>3%</div> <div> <div></div> <div>67%</div> <div>17%</div> <div>•</div> <div>15%</div> </div> </div>
1	C	255	<div> <div>4%</div> <div> <div></div> <div>66%</div> <div>18%</div> <div>•</div> <div>13%</div> </div> </div>
1	D	255	<div> <div>2%</div> <div> <div></div> <div>69%</div> <div>16%</div> <div>•</div> <div>14%</div> </div> </div>
1	E	255	<div> <div>4%</div> <div> <div></div> <div>67%</div> <div>17%</div> <div>•</div> <div>14%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	255	<div><div><div></div><div></div><div></div></div><div><div>2%</div><div>67%</div><div>18%</div><div>•</div><div>14%</div></div></div>
1	G	255	<div><div><div></div><div></div><div></div></div><div><div>2%</div><div>71%</div><div>13%</div><div>•</div><div>15%</div></div></div>
1	H	255	<div><div><div></div><div></div><div></div></div><div><div>3%</div><div>66%</div><div>18%</div><div>•</div><div>15%</div></div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14128 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 Fluorescent protein Dronpa.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	0	0
			1726	1105	292	320	9			
1	B	216	Total	C	N	O	S	0	0	0
			1730	1108	293	320	9			
1	C	221	Total	C	N	O	S	0	0	0
			1743	1116	295	323	9			
1	D	220	Total	C	N	O	S	0	0	0
			1741	1113	293	326	9			
1	E	219	Total	C	N	O	S	0	1	0
			1745	1117	298	321	9			
1	F	219	Total	C	N	O	S	0	0	0
			1756	1124	296	327	9			
1	G	216	Total	C	N	O	S	0	0	0
			1720	1101	288	322	9			
1	H	216	Total	C	N	O	S	0	0	0
			1710	1093	288	320	9			

There are 328 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-27	GLY	-	expression tag	UNP Q5TLG6
A	-26	SER	-	expression tag	UNP Q5TLG6
A	-25	SER	-	expression tag	UNP Q5TLG6
A	-24	HIS	-	expression tag	UNP Q5TLG6
A	-23	HIS	-	expression tag	UNP Q5TLG6
A	-22	HIS	-	expression tag	UNP Q5TLG6
A	-21	HIS	-	expression tag	UNP Q5TLG6
A	-20	HIS	-	expression tag	UNP Q5TLG6
A	-19	HIS	-	expression tag	UNP Q5TLG6
A	-18	SER	-	expression tag	UNP Q5TLG6
A	-17	SER	-	expression tag	UNP Q5TLG6
A	-16	GLY	-	expression tag	UNP Q5TLG6
A	-15	LEU	-	expression tag	UNP Q5TLG6

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	VAL	-	expression tag	UNP Q5TLG6
A	-13	PRO	-	expression tag	UNP Q5TLG6
A	-12	GLY	-	expression tag	UNP Q5TLG6
A	-11	GLY	-	expression tag	UNP Q5TLG6
A	-10	SER	-	expression tag	UNP Q5TLG6
A	-9	HIS	-	expression tag	UNP Q5TLG6
A	-8	MET	-	expression tag	UNP Q5TLG6
A	-7	VAL	-	expression tag	UNP Q5TLG6
A	-6	SER	-	expression tag	UNP Q5TLG6
A	-5	LYS	-	expression tag	UNP Q5TLG6
A	-4	GLY	-	expression tag	UNP Q5TLG6
A	-3	GLU	-	expression tag	UNP Q5TLG6
A	-2	GLU	-	expression tag	UNP Q5TLG6
A	-1	ASN	-	expression tag	UNP Q5TLG6
A	0	ASN	-	expression tag	UNP Q5TLG6
A	1	MET	-	expression tag	UNP Q5TLG6
A	2	ALA	-	expression tag	UNP Q5TLG6
A	63	GYC	CYS	chromophore	UNP Q5TLG6
A	63	GYC	TYR	chromophore	UNP Q5TLG6
A	63	GYC	GLY	chromophore	UNP Q5TLG6
A	142	ALA	SER	engineered mutation	UNP Q5TLG6
A	159	THR	MET	engineered mutation	UNP Q5TLG6
A	218	GLY	GLU	conflict	UNP Q5TLG6
A	224	MET	-	insertion	UNP Q5TLG6
A	225	ASP	-	insertion	UNP Q5TLG6
A	226	GLU	-	insertion	UNP Q5TLG6
A	227	LEU	-	insertion	UNP Q5TLG6
A	228	TYR	-	insertion	UNP Q5TLG6
B	-27	GLY	-	expression tag	UNP Q5TLG6
B	-26	SER	-	expression tag	UNP Q5TLG6
B	-25	SER	-	expression tag	UNP Q5TLG6
B	-24	HIS	-	expression tag	UNP Q5TLG6
B	-23	HIS	-	expression tag	UNP Q5TLG6
B	-22	HIS	-	expression tag	UNP Q5TLG6
B	-21	HIS	-	expression tag	UNP Q5TLG6
B	-20	HIS	-	expression tag	UNP Q5TLG6
B	-19	HIS	-	expression tag	UNP Q5TLG6
B	-18	SER	-	expression tag	UNP Q5TLG6
B	-17	SER	-	expression tag	UNP Q5TLG6
B	-16	GLY	-	expression tag	UNP Q5TLG6
B	-15	LEU	-	expression tag	UNP Q5TLG6
B	-14	VAL	-	expression tag	UNP Q5TLG6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-13	PRO	-	expression tag	UNP Q5TLG6
B	-12	GLY	-	expression tag	UNP Q5TLG6
B	-11	GLY	-	expression tag	UNP Q5TLG6
B	-10	SER	-	expression tag	UNP Q5TLG6
B	-9	HIS	-	expression tag	UNP Q5TLG6
B	-8	MET	-	expression tag	UNP Q5TLG6
B	-7	VAL	-	expression tag	UNP Q5TLG6
B	-6	SER	-	expression tag	UNP Q5TLG6
B	-5	LYS	-	expression tag	UNP Q5TLG6
B	-4	GLY	-	expression tag	UNP Q5TLG6
B	-3	GLU	-	expression tag	UNP Q5TLG6
B	-2	GLU	-	expression tag	UNP Q5TLG6
B	-1	ASN	-	expression tag	UNP Q5TLG6
B	0	ASN	-	expression tag	UNP Q5TLG6
B	1	MET	-	expression tag	UNP Q5TLG6
B	2	ALA	-	expression tag	UNP Q5TLG6
B	63	GYC	CYS	chromophore	UNP Q5TLG6
B	63	GYC	TYR	chromophore	UNP Q5TLG6
B	63	GYC	GLY	chromophore	UNP Q5TLG6
B	142	ALA	SER	engineered mutation	UNP Q5TLG6
B	159	THR	MET	engineered mutation	UNP Q5TLG6
B	218	GLY	GLU	conflict	UNP Q5TLG6
B	224	MET	-	insertion	UNP Q5TLG6
B	225	ASP	-	insertion	UNP Q5TLG6
B	226	GLU	-	insertion	UNP Q5TLG6
B	227	LEU	-	insertion	UNP Q5TLG6
B	228	TYR	-	insertion	UNP Q5TLG6
C	-27	GLY	-	expression tag	UNP Q5TLG6
C	-26	SER	-	expression tag	UNP Q5TLG6
C	-25	SER	-	expression tag	UNP Q5TLG6
C	-24	HIS	-	expression tag	UNP Q5TLG6
C	-23	HIS	-	expression tag	UNP Q5TLG6
C	-22	HIS	-	expression tag	UNP Q5TLG6
C	-21	HIS	-	expression tag	UNP Q5TLG6
C	-20	HIS	-	expression tag	UNP Q5TLG6
C	-19	HIS	-	expression tag	UNP Q5TLG6
C	-18	SER	-	expression tag	UNP Q5TLG6
C	-17	SER	-	expression tag	UNP Q5TLG6
C	-16	GLY	-	expression tag	UNP Q5TLG6
C	-15	LEU	-	expression tag	UNP Q5TLG6
C	-14	VAL	-	expression tag	UNP Q5TLG6
C	-13	PRO	-	expression tag	UNP Q5TLG6

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-12	GLY	-	expression tag	UNP Q5TLG6
C	-11	GLY	-	expression tag	UNP Q5TLG6
C	-10	SER	-	expression tag	UNP Q5TLG6
C	-9	HIS	-	expression tag	UNP Q5TLG6
C	-8	MET	-	expression tag	UNP Q5TLG6
C	-7	VAL	-	expression tag	UNP Q5TLG6
C	-6	SER	-	expression tag	UNP Q5TLG6
C	-5	LYS	-	expression tag	UNP Q5TLG6
C	-4	GLY	-	expression tag	UNP Q5TLG6
C	-3	GLU	-	expression tag	UNP Q5TLG6
C	-2	GLU	-	expression tag	UNP Q5TLG6
C	-1	ASN	-	expression tag	UNP Q5TLG6
C	0	ASN	-	expression tag	UNP Q5TLG6
C	1	MET	-	expression tag	UNP Q5TLG6
C	2	ALA	-	expression tag	UNP Q5TLG6
C	63	GYC	CYS	chromophore	UNP Q5TLG6
C	63	GYC	TYR	chromophore	UNP Q5TLG6
C	63	GYC	GLY	chromophore	UNP Q5TLG6
C	142	ALA	SER	engineered mutation	UNP Q5TLG6
C	159	THR	MET	engineered mutation	UNP Q5TLG6
C	218	GLY	GLU	conflict	UNP Q5TLG6
C	224	MET	-	insertion	UNP Q5TLG6
C	225	ASP	-	insertion	UNP Q5TLG6
C	226	GLU	-	insertion	UNP Q5TLG6
C	227	LEU	-	insertion	UNP Q5TLG6
C	228	TYR	-	insertion	UNP Q5TLG6
D	-27	GLY	-	expression tag	UNP Q5TLG6
D	-26	SER	-	expression tag	UNP Q5TLG6
D	-25	SER	-	expression tag	UNP Q5TLG6
D	-24	HIS	-	expression tag	UNP Q5TLG6
D	-23	HIS	-	expression tag	UNP Q5TLG6
D	-22	HIS	-	expression tag	UNP Q5TLG6
D	-21	HIS	-	expression tag	UNP Q5TLG6
D	-20	HIS	-	expression tag	UNP Q5TLG6
D	-19	HIS	-	expression tag	UNP Q5TLG6
D	-18	SER	-	expression tag	UNP Q5TLG6
D	-17	SER	-	expression tag	UNP Q5TLG6
D	-16	GLY	-	expression tag	UNP Q5TLG6
D	-15	LEU	-	expression tag	UNP Q5TLG6
D	-14	VAL	-	expression tag	UNP Q5TLG6
D	-13	PRO	-	expression tag	UNP Q5TLG6
D	-12	GLY	-	expression tag	UNP Q5TLG6

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-11	GLY	-	expression tag	UNP Q5TLG6
D	-10	SER	-	expression tag	UNP Q5TLG6
D	-9	HIS	-	expression tag	UNP Q5TLG6
D	-8	MET	-	expression tag	UNP Q5TLG6
D	-7	VAL	-	expression tag	UNP Q5TLG6
D	-6	SER	-	expression tag	UNP Q5TLG6
D	-5	LYS	-	expression tag	UNP Q5TLG6
D	-4	GLY	-	expression tag	UNP Q5TLG6
D	-3	GLU	-	expression tag	UNP Q5TLG6
D	-2	GLU	-	expression tag	UNP Q5TLG6
D	-1	ASN	-	expression tag	UNP Q5TLG6
D	0	ASN	-	expression tag	UNP Q5TLG6
D	1	MET	-	expression tag	UNP Q5TLG6
D	2	ALA	-	expression tag	UNP Q5TLG6
D	63	GYC	CYS	chromophore	UNP Q5TLG6
D	63	GYC	TYR	chromophore	UNP Q5TLG6
D	63	GYC	GLY	chromophore	UNP Q5TLG6
D	142	ALA	SER	engineered mutation	UNP Q5TLG6
D	159	THR	MET	engineered mutation	UNP Q5TLG6
D	218	GLY	GLU	conflict	UNP Q5TLG6
D	224	MET	-	insertion	UNP Q5TLG6
D	225	ASP	-	insertion	UNP Q5TLG6
D	226	GLU	-	insertion	UNP Q5TLG6
D	227	LEU	-	insertion	UNP Q5TLG6
D	228	TYR	-	insertion	UNP Q5TLG6
E	-27	GLY	-	expression tag	UNP Q5TLG6
E	-26	SER	-	expression tag	UNP Q5TLG6
E	-25	SER	-	expression tag	UNP Q5TLG6
E	-24	HIS	-	expression tag	UNP Q5TLG6
E	-23	HIS	-	expression tag	UNP Q5TLG6
E	-22	HIS	-	expression tag	UNP Q5TLG6
E	-21	HIS	-	expression tag	UNP Q5TLG6
E	-20	HIS	-	expression tag	UNP Q5TLG6
E	-19	HIS	-	expression tag	UNP Q5TLG6
E	-18	SER	-	expression tag	UNP Q5TLG6
E	-17	SER	-	expression tag	UNP Q5TLG6
E	-16	GLY	-	expression tag	UNP Q5TLG6
E	-15	LEU	-	expression tag	UNP Q5TLG6
E	-14	VAL	-	expression tag	UNP Q5TLG6
E	-13	PRO	-	expression tag	UNP Q5TLG6
E	-12	GLY	-	expression tag	UNP Q5TLG6
E	-11	GLY	-	expression tag	UNP Q5TLG6

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-10	SER	-	expression tag	UNP Q5TLG6
E	-9	HIS	-	expression tag	UNP Q5TLG6
E	-8	MET	-	expression tag	UNP Q5TLG6
E	-7	VAL	-	expression tag	UNP Q5TLG6
E	-6	SER	-	expression tag	UNP Q5TLG6
E	-5	LYS	-	expression tag	UNP Q5TLG6
E	-4	GLY	-	expression tag	UNP Q5TLG6
E	-3	GLU	-	expression tag	UNP Q5TLG6
E	-2	GLU	-	expression tag	UNP Q5TLG6
E	-1	ASN	-	expression tag	UNP Q5TLG6
E	0	ASN	-	expression tag	UNP Q5TLG6
E	1	MET	-	expression tag	UNP Q5TLG6
E	2	ALA	-	expression tag	UNP Q5TLG6
E	63	GYC	CYS	chromophore	UNP Q5TLG6
E	63	GYC	TYR	chromophore	UNP Q5TLG6
E	63	GYC	GLY	chromophore	UNP Q5TLG6
E	142	ALA	SER	engineered mutation	UNP Q5TLG6
E	159	THR	MET	engineered mutation	UNP Q5TLG6
E	218	GLY	GLU	conflict	UNP Q5TLG6
E	224	MET	-	insertion	UNP Q5TLG6
E	225	ASP	-	insertion	UNP Q5TLG6
E	226	GLU	-	insertion	UNP Q5TLG6
E	227	LEU	-	insertion	UNP Q5TLG6
E	228	TYR	-	insertion	UNP Q5TLG6
F	-27	GLY	-	expression tag	UNP Q5TLG6
F	-26	SER	-	expression tag	UNP Q5TLG6
F	-25	SER	-	expression tag	UNP Q5TLG6
F	-24	HIS	-	expression tag	UNP Q5TLG6
F	-23	HIS	-	expression tag	UNP Q5TLG6
F	-22	HIS	-	expression tag	UNP Q5TLG6
F	-21	HIS	-	expression tag	UNP Q5TLG6
F	-20	HIS	-	expression tag	UNP Q5TLG6
F	-19	HIS	-	expression tag	UNP Q5TLG6
F	-18	SER	-	expression tag	UNP Q5TLG6
F	-17	SER	-	expression tag	UNP Q5TLG6
F	-16	GLY	-	expression tag	UNP Q5TLG6
F	-15	LEU	-	expression tag	UNP Q5TLG6
F	-14	VAL	-	expression tag	UNP Q5TLG6
F	-13	PRO	-	expression tag	UNP Q5TLG6
F	-12	GLY	-	expression tag	UNP Q5TLG6
F	-11	GLY	-	expression tag	UNP Q5TLG6
F	-10	SER	-	expression tag	UNP Q5TLG6

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-9	HIS	-	expression tag	UNP Q5TLG6
F	-8	MET	-	expression tag	UNP Q5TLG6
F	-7	VAL	-	expression tag	UNP Q5TLG6
F	-6	SER	-	expression tag	UNP Q5TLG6
F	-5	LYS	-	expression tag	UNP Q5TLG6
F	-4	GLY	-	expression tag	UNP Q5TLG6
F	-3	GLU	-	expression tag	UNP Q5TLG6
F	-2	GLU	-	expression tag	UNP Q5TLG6
F	-1	ASN	-	expression tag	UNP Q5TLG6
F	0	ASN	-	expression tag	UNP Q5TLG6
F	1	MET	-	expression tag	UNP Q5TLG6
F	2	ALA	-	expression tag	UNP Q5TLG6
F	63	GYC	CYS	chromophore	UNP Q5TLG6
F	63	GYC	TYR	chromophore	UNP Q5TLG6
F	63	GYC	GLY	chromophore	UNP Q5TLG6
F	142	ALA	SER	engineered mutation	UNP Q5TLG6
F	159	THR	MET	engineered mutation	UNP Q5TLG6
F	218	GLY	GLU	conflict	UNP Q5TLG6
F	224	MET	-	insertion	UNP Q5TLG6
F	225	ASP	-	insertion	UNP Q5TLG6
F	226	GLU	-	insertion	UNP Q5TLG6
F	227	LEU	-	insertion	UNP Q5TLG6
F	228	TYR	-	insertion	UNP Q5TLG6
G	-27	GLY	-	expression tag	UNP Q5TLG6
G	-26	SER	-	expression tag	UNP Q5TLG6
G	-25	SER	-	expression tag	UNP Q5TLG6
G	-24	HIS	-	expression tag	UNP Q5TLG6
G	-23	HIS	-	expression tag	UNP Q5TLG6
G	-22	HIS	-	expression tag	UNP Q5TLG6
G	-21	HIS	-	expression tag	UNP Q5TLG6
G	-20	HIS	-	expression tag	UNP Q5TLG6
G	-19	HIS	-	expression tag	UNP Q5TLG6
G	-18	SER	-	expression tag	UNP Q5TLG6
G	-17	SER	-	expression tag	UNP Q5TLG6
G	-16	GLY	-	expression tag	UNP Q5TLG6
G	-15	LEU	-	expression tag	UNP Q5TLG6
G	-14	VAL	-	expression tag	UNP Q5TLG6
G	-13	PRO	-	expression tag	UNP Q5TLG6
G	-12	GLY	-	expression tag	UNP Q5TLG6
G	-11	GLY	-	expression tag	UNP Q5TLG6
G	-10	SER	-	expression tag	UNP Q5TLG6
G	-9	HIS	-	expression tag	UNP Q5TLG6

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-8	MET	-	expression tag	UNP Q5TLG6
G	-7	VAL	-	expression tag	UNP Q5TLG6
G	-6	SER	-	expression tag	UNP Q5TLG6
G	-5	LYS	-	expression tag	UNP Q5TLG6
G	-4	GLY	-	expression tag	UNP Q5TLG6
G	-3	GLU	-	expression tag	UNP Q5TLG6
G	-2	GLU	-	expression tag	UNP Q5TLG6
G	-1	ASN	-	expression tag	UNP Q5TLG6
G	0	ASN	-	expression tag	UNP Q5TLG6
G	1	MET	-	expression tag	UNP Q5TLG6
G	2	ALA	-	expression tag	UNP Q5TLG6
G	63	GYC	CYS	chromophore	UNP Q5TLG6
G	63	GYC	TYR	chromophore	UNP Q5TLG6
G	63	GYC	GLY	chromophore	UNP Q5TLG6
G	142	ALA	SER	engineered mutation	UNP Q5TLG6
G	159	THR	MET	engineered mutation	UNP Q5TLG6
G	218	GLY	GLU	conflict	UNP Q5TLG6
G	224	MET	-	insertion	UNP Q5TLG6
G	225	ASP	-	insertion	UNP Q5TLG6
G	226	GLU	-	insertion	UNP Q5TLG6
G	227	LEU	-	insertion	UNP Q5TLG6
G	228	TYR	-	insertion	UNP Q5TLG6
H	-27	GLY	-	expression tag	UNP Q5TLG6
H	-26	SER	-	expression tag	UNP Q5TLG6
H	-25	SER	-	expression tag	UNP Q5TLG6
H	-24	HIS	-	expression tag	UNP Q5TLG6
H	-23	HIS	-	expression tag	UNP Q5TLG6
H	-22	HIS	-	expression tag	UNP Q5TLG6
H	-21	HIS	-	expression tag	UNP Q5TLG6
H	-20	HIS	-	expression tag	UNP Q5TLG6
H	-19	HIS	-	expression tag	UNP Q5TLG6
H	-18	SER	-	expression tag	UNP Q5TLG6
H	-17	SER	-	expression tag	UNP Q5TLG6
H	-16	GLY	-	expression tag	UNP Q5TLG6
H	-15	LEU	-	expression tag	UNP Q5TLG6
H	-14	VAL	-	expression tag	UNP Q5TLG6
H	-13	PRO	-	expression tag	UNP Q5TLG6
H	-12	GLY	-	expression tag	UNP Q5TLG6
H	-11	GLY	-	expression tag	UNP Q5TLG6
H	-10	SER	-	expression tag	UNP Q5TLG6
H	-9	HIS	-	expression tag	UNP Q5TLG6
H	-8	MET	-	expression tag	UNP Q5TLG6

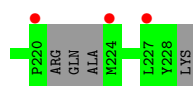
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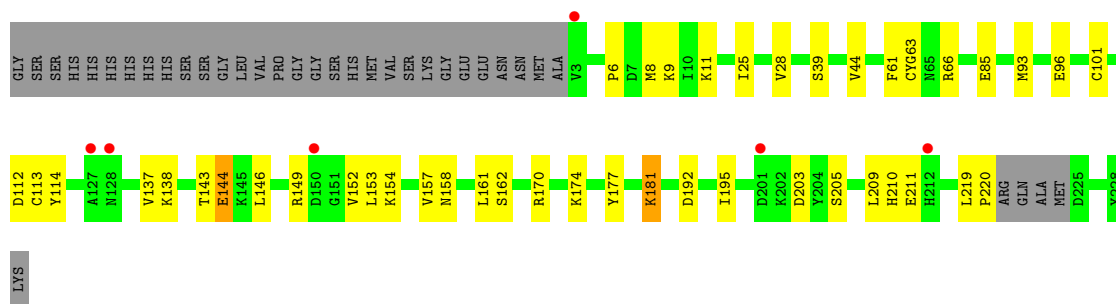
Chain	Residue	Modelled	Actual	Comment	Reference
H	-7	VAL	-	expression tag	UNP Q5TLG6
H	-6	SER	-	expression tag	UNP Q5TLG6
H	-5	LYS	-	expression tag	UNP Q5TLG6
H	-4	GLY	-	expression tag	UNP Q5TLG6
H	-3	GLU	-	expression tag	UNP Q5TLG6
H	-2	GLU	-	expression tag	UNP Q5TLG6
H	-1	ASN	-	expression tag	UNP Q5TLG6
H	0	ASN	-	expression tag	UNP Q5TLG6
H	1	MET	-	expression tag	UNP Q5TLG6
H	2	ALA	-	expression tag	UNP Q5TLG6
H	63	GYC	CYS	chromophore	UNP Q5TLG6
H	63	GYC	TYR	chromophore	UNP Q5TLG6
H	63	GYC	GLY	chromophore	UNP Q5TLG6
H	142	ALA	SER	engineered mutation	UNP Q5TLG6
H	159	THR	MET	engineered mutation	UNP Q5TLG6
H	218	GLY	GLU	conflict	UNP Q5TLG6
H	224	MET	-	insertion	UNP Q5TLG6
H	225	ASP	-	insertion	UNP Q5TLG6
H	226	GLU	-	insertion	UNP Q5TLG6
H	227	LEU	-	insertion	UNP Q5TLG6
H	228	TYR	-	insertion	UNP Q5TLG6

- Molecule 2 is water.

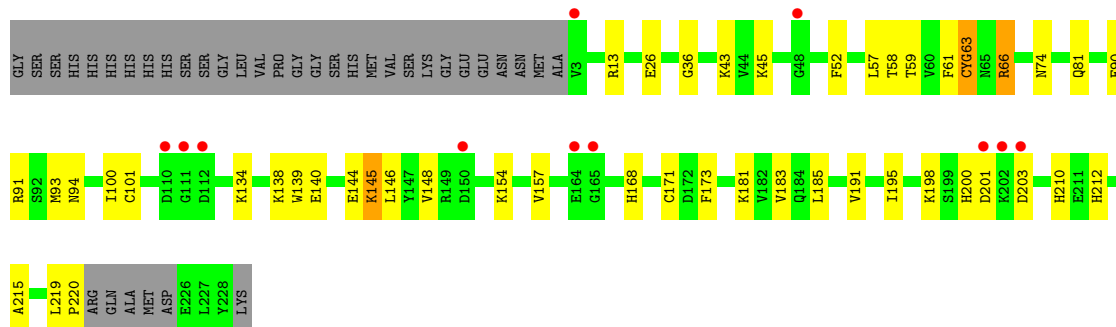
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	38	Total O 38 38	0	0
2	B	37	Total O 37 37	0	0
2	C	25	Total O 25 25	0	0
2	D	31	Total O 31 31	0	0
2	E	29	Total O 29 29	0	0
2	F	31	Total O 31 31	0	0
2	G	30	Total O 30 30	0	0
2	H	36	Total O 36 36	0	0



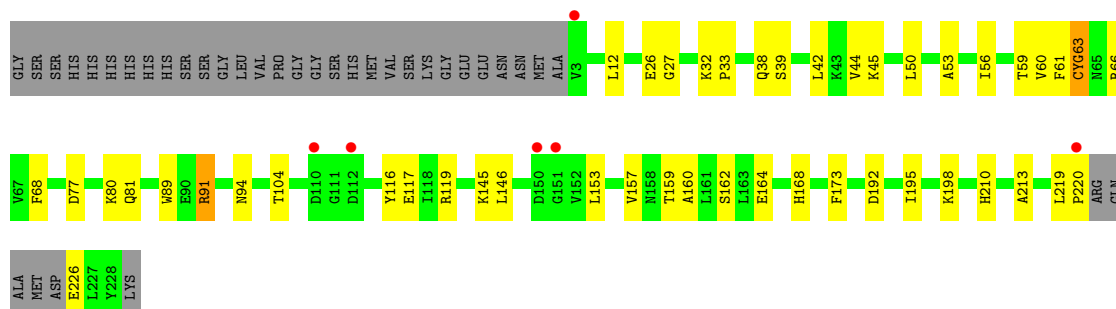
• Molecule 1: Fluorescent protein Dronpa



• Molecule 1: Fluorescent protein Dronpa

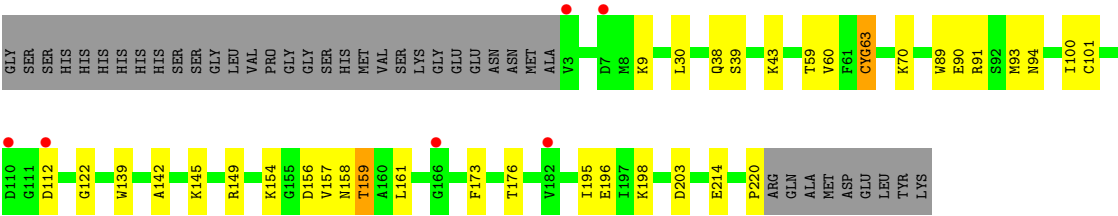


• Molecule 1: Fluorescent protein Dronpa

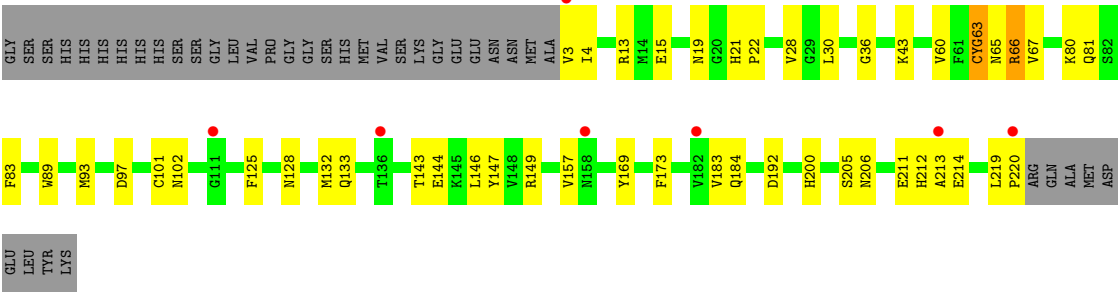


• Molecule 1: Fluorescent protein Dronpa





• Molecule 1: Fluorescent protein Dronpa



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.41Å 85.81Å 143.17Å 90.00° 95.01° 90.00°	Depositor
Resolution (Å)	37.29 – 2.64 39.00 – 2.64	Depositor EDS
% Data completeness (in resolution range)	94.6 (37.29-2.64) 82.7 (39.00-2.64)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.12 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.13rc2_2986	Depositor
R, R_{free}	0.226 , 0.284 0.226 , 0.275	Depositor DCC
R_{free} test set	2674 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.472	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 40.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	14128	wwPDB-VP
Average B, all atoms (Å ²)	38.0	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 48.98 % 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 7.9617e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

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

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	0.65	0/1751	0.66	0/2368
1	B	0.65	0/1755	0.66	0/2372
1	C	0.64	0/1767	0.69	0/2391
1	D	0.68	0/1765	0.69	0/2389
1	E	0.71	0/1769	0.69	0/2390
1	F	0.63	0/1780	0.65	0/2403
1	G	0.69	0/1745	0.70	0/2362
1	H	0.69	0/1735	0.68	0/2352
All	All	0.67	0/14067	0.68	0/19027

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	F	0	1
All	All	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	61	PHE	Mainchain
1	E	61	PHE	Mainchain

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Mol	Chain	Res	Type	Group
1	F	61	PHE	Mainchain

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	1726	0	1650	28	0
1	B	1730	0	1661	29	0
1	C	1743	0	1646	42	0
1	D	1741	0	1638	30	0
1	E	1745	0	1664	41	0
1	F	1756	0	1684	28	0
1	G	1720	0	1632	33	0
1	H	1710	0	1605	45	0
2	A	38	0	0	4	0
2	B	37	0	0	2	0
2	C	25	0	0	0	0
2	D	31	0	0	2	0
2	E	29	0	0	2	0
2	F	31	0	0	2	0
2	G	30	0	0	2	0
2	H	36	0	0	2	0
All	All	14128	0	13180	254	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 9.

All (254) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:144:GLU:HG3	1:H:157:VAL:HG23	1.32	1.09
1:C:31:GLY:HA3	1:C:68:PHE:CE1	1.88	1.07
1:H:144:GLU:HG3	1:H:157:VAL:CG2	1.91	0.98
1:E:157:VAL:HG22	1:E:173:PHE:HB2	1.55	0.88
1:E:58:THR:O	1:E:63:GYC:HB11	1.74	0.87
1:E:138:LYS:HE3	1:E:139:TRP:O	1.76	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:138:LYS:HG2	1:E:139:TRP:N	1.90	0.85
1:G:94:ASN:OD1	1:G:100:ILE:HD13	1.77	0.84
1:A:9:LYS:HD3	1:A:30:LEU:HB3	1.57	0.84
1:H:19:ASN:HD21	1:H:125:PHE:HB2	1.44	0.83
1:G:139:TRP:CD1	1:G:161:LEU:HD21	2.14	0.82
1:G:139:TRP:NE1	1:G:161:LEU:HD21	1.95	0.79
1:G:59:THR:HG21	1:G:173:PHE:HE1	1.51	0.75
1:C:27:GLY:HA3	1:C:42:LEU:HD12	1.66	0.75
1:E:168:HIS:ND1	2:E:301:HOH:O	2.21	0.73
1:C:168:HIS:O	1:G:149:ARG:NH2	2.22	0.72
1:B:157:VAL:HB	1:B:173:PHE:HB2	1.71	0.71
1:F:159:THR:HG22	1:F:160:ALA:H	1.56	0.71
1:H:67:VAL:HG11	1:H:83:PHE:HE2	1.56	0.70
1:C:33:PRO:O	1:C:80:LYS:CE	2.42	0.67
1:D:11:LYS:HG3	1:D:28:VAL:HG12	1.77	0.67
1:C:148:VAL:HG21	1:C:185:LEU:HB3	1.77	0.67
1:A:3:VAL:N	2:A:302:HOH:O	2.27	0.67
1:A:5:LYS:HB2	1:A:8:MET:SD	2.35	0.67
1:C:5:LYS:HG2	1:C:7:ASP:H	1.61	0.66
1:B:76:VAL:HG11	1:B:184:GLN:HG2	1.77	0.66
1:D:195:ILE:HD11	1:D:209:LEU:HD21	1.77	0.65
1:E:52:PHE:HE1	1:E:57:LEU:HD11	1.59	0.65
1:E:144:GLU:HB2	1:E:157:VAL:HG12	1.76	0.65
1:C:30:LEU:HD12	1:C:30:LEU:O	1.96	0.65
1:E:66:ARG:HG2	1:E:66:ARG:HH11	1.62	0.64
1:E:157:VAL:HG22	1:E:173:PHE:CB	2.27	0.64
1:G:157:VAL:HB	1:G:173:PHE:HB2	1.80	0.64
1:A:149:ARG:NH2	1:E:168:HIS:O	2.31	0.63
1:B:100:ILE:HG22	1:H:102:ASN:HD21	1.62	0.63
1:B:145:LYS:HG2	1:B:156:ASP:HB2	1.80	0.63
1:H:3:VAL:HG13	1:H:4:ILE:HG13	1.81	0.63
1:G:154:LYS:NZ	1:G:176:THR:OG1	2.27	0.62
1:B:40:MET:HE1	1:B:61:PHE:HB3	1.81	0.62
1:H:67:VAL:O	1:H:80:LYS:NZ	2.32	0.62
1:E:59:THR:O	1:E:91:ARG:NH2	2.32	0.62
1:H:144:GLU:CG	1:H:157:VAL:CG2	2.73	0.62
1:C:33:PRO:O	1:C:80:LYS:NZ	2.33	0.60
1:E:26:GLU:HG3	1:E:45:LYS:HD2	1.84	0.60
1:F:26:GLU:HB2	1:F:45:LYS:HE3	1.84	0.60
1:B:149:ARG:NH2	1:F:168:HIS:O	2.35	0.59
1:F:59:THR:HG21	1:F:173:PHE:CE2	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:ASP:O	1:B:81:GLN:HG3	2.03	0.59
1:C:31:GLY:HA3	1:C:68:PHE:CZ	2.36	0.59
1:D:219:LEU:HD21	1:H:214:GLU:HG3	1.85	0.59
1:G:203:ASP:OD1	1:G:203:ASP:N	2.28	0.59
1:C:139:TRP:HB2	1:C:195:ILE:CG2	2.32	0.59
1:G:93:MET:HB2	1:G:101:CYS:HB2	1.84	0.59
1:H:219:LEU:N	1:H:219:LEU:HD12	2.18	0.58
1:B:158:ASN:O	1:F:145:LYS:CD	2.52	0.58
1:C:139:TRP:CD2	1:C:161:LEU:HD13	2.39	0.58
1:E:144:GLU:CB	1:E:157:VAL:HG12	2.34	0.57
1:B:26:GLU:OE2	1:B:45:LYS:HE2	2.05	0.57
1:D:195:ILE:HG22	2:D:304:HOH:O	2.04	0.57
1:C:30:LEU:HD12	1:C:39:SER:HB3	1.87	0.57
1:F:44:VAL:HG21	1:F:50:LEU:HD21	1.87	0.56
1:G:139:TRP:CD1	1:G:161:LEU:CD2	2.88	0.56
1:C:60:VAL:HG13	1:C:89:TRP:HH2	1.69	0.56
1:E:144:GLU:HB2	1:E:157:VAL:CG1	2.35	0.56
1:D:192:ASP:HB3	1:H:219:LEU:HD11	1.87	0.56
1:F:12:LEU:HB2	1:F:116:TYR:HB2	1.87	0.56
1:E:157:VAL:HG22	1:E:173:PHE:CG	2.40	0.56
1:G:70:LYS:HB3	1:G:214:GLU:HG2	1.88	0.55
1:C:31:GLY:HA3	1:C:68:PHE:HE1	1.63	0.55
1:B:124:ASN:HB3	2:B:316:HOH:O	2.06	0.55
1:D:170:ARG:NH1	1:H:147:TYR:OH	2.40	0.55
1:G:94:ASN:OD1	1:G:100:ILE:CD1	2.54	0.55
1:B:200:HIS:ND1	1:B:201:ASP:O	2.40	0.55
1:C:27:GLY:HA3	1:C:42:LEU:CD1	2.36	0.55
1:B:200:HIS:NE2	2:B:304:HOH:O	2.33	0.55
1:E:200:HIS:ND1	1:E:201:ASP:O	2.39	0.55
1:G:59:THR:HG22	1:G:91:ARG:HH12	1.73	0.55
1:D:9:LYS:HE3	1:D:112:ASP:OD2	2.06	0.54
1:E:146:LEU:N	1:E:146:LEU:HD12	2.21	0.54
1:C:85:GLU:HG2	1:C:181:LYS:HE2	1.88	0.54
1:B:7:ASP:OD1	1:B:32:LYS:NZ	2.40	0.54
1:H:67:VAL:HG11	1:H:83:PHE:CE2	2.40	0.54
1:A:154:LYS:HG2	2:A:333:HOH:O	2.07	0.54
1:D:195:ILE:HD12	1:D:211:GLU:HG3	1.89	0.54
1:H:144:GLU:CB	1:H:157:VAL:HG22	2.37	0.54
1:B:60:VAL:HG13	1:B:89:TRP:HH2	1.72	0.54
1:D:143:THR:HB	1:H:143:THR:HG21	1.90	0.54
1:F:60:VAL:HG13	1:F:89:TRP:HH2	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:144:GLU:HB2	1:H:157:VAL:HG22	1.90	0.54
1:C:159:THR:HG22	1:C:160:ALA:H	1.73	0.53
1:A:63:GYC:O	1:A:89:TRP:NE1	2.36	0.53
1:A:157:VAL:HB	1:A:173:PHE:HB2	1.91	0.53
1:C:30:LEU:CD1	1:C:39:SER:HB3	2.39	0.53
1:G:139:TRP:CE2	1:G:161:LEU:HG	2.44	0.53
1:D:66:ARG:NH2	1:D:177:TYR:OH	2.39	0.53
1:B:192:ASP:O	1:B:213:ALA:HA	2.08	0.53
1:H:60:VAL:HG13	1:H:89:TRP:HH2	1.74	0.53
1:C:19:ASN:HA	1:E:90:GLU:OE1	2.10	0.52
1:H:43:LYS:HA	1:H:205:SER:O	2.09	0.52
1:C:146:LEU:N	1:C:146:LEU:HD12	2.24	0.52
1:D:137:VAL:HB	1:D:162:SER:OG	2.10	0.52
1:F:77:ASP:O	1:F:81:GLN:HG3	2.09	0.52
1:C:10:ILE:N	1:C:29:GLY:O	2.38	0.52
1:A:128:ASN:HA	1:A:133:GLN:HE21	1.75	0.52
1:G:30:LEU:HD23	1:G:30:LEU:N	2.24	0.52
1:H:219:LEU:N	1:H:219:LEU:CD1	2.73	0.52
1:A:159:THR:HA	2:A:317:HOH:O	2.11	0.51
1:C:137:VAL:HB	1:C:162:SER:OG	2.11	0.51
1:F:226:GLU:N	2:F:307:HOH:O	2.43	0.51
1:E:94[A]:ASN:ND2	2:E:306:HOH:O	2.43	0.51
1:C:157:VAL:HB	1:C:173:PHE:HB2	1.92	0.51
1:A:60:VAL:HG13	1:A:89:TRP:HH2	1.75	0.50
1:D:144:GLU:HB2	1:D:157:VAL:HG22	1.92	0.50
1:H:65:ASN:CG	1:H:67:VAL:HG22	2.32	0.50
1:F:104:THR:HG22	1:F:119:ARG:HB3	1.92	0.50
1:B:158:ASN:O	1:F:145:LYS:HD3	2.11	0.50
1:C:160:ALA:HB1	1:C:168:HIS:HB3	1.93	0.50
1:D:8:MET:HG3	1:D:114:TYR:HE2	1.77	0.49
1:C:27:GLY:CA	1:C:42:LEU:HD12	2.39	0.49
1:A:43:LYS:NZ	2:A:307:HOH:O	2.44	0.49
1:C:36:GLY:O	1:C:212:HIS:HA	2.13	0.49
1:G:139:TRP:CZ2	1:G:161:LEU:HD11	2.48	0.49
1:H:183:VAL:O	2:H:301:HOH:O	2.20	0.49
1:E:144:GLU:HA	1:E:157:VAL:HG12	1.94	0.49
1:C:194:HIS:CG	1:G:220:PRO:HG3	2.47	0.49
1:B:66:ARG:HD2	1:B:191:VAL:HG11	1.95	0.49
1:F:63:GYC:HB11	1:F:195:ILE:HD12	1.95	0.49
1:A:100:ILE:HD12	2:G:325:HOH:O	2.12	0.49
1:E:148:VAL:HG21	1:E:185:LEU:HB3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:96:GLU:CD	1:H:149:ARG:HH22	2.16	0.48
1:G:90:GLU:HG2	1:G:176:THR:HB	1.95	0.48
1:H:81:GLN:HB3	1:H:183:VAL:CG1	2.44	0.48
1:E:13:ARG:NH1	1:E:26:GLU:OE2	2.47	0.48
1:E:171:CYS:HG	1:E:173:PHE:HE1	1.57	0.48
1:C:5:LYS:HG3	1:C:6:PRO:HD2	1.96	0.47
1:E:203:ASP:OD2	1:E:203:ASP:N	2.42	0.47
1:D:138:LYS:O	1:D:161:LEU:HD23	2.14	0.47
1:D:219:LEU:HD12	1:D:220:PRO:HD2	1.95	0.47
1:B:137:VAL:HG23	1:B:164:GLU:N	2.30	0.47
1:C:31:GLY:CA	1:C:68:PHE:CE1	2.80	0.47
1:H:183:VAL:HG12	1:H:184:GLN:N	2.29	0.47
1:F:104:THR:HG23	2:F:305:HOH:O	2.15	0.47
1:H:4:ILE:HB	2:H:309:HOH:O	2.14	0.47
1:C:163:LEU:HD11	1:C:169:TYR:HB2	1.96	0.47
1:C:194:HIS:CD2	1:G:220:PRO:HG3	2.50	0.47
1:E:66:ARG:HG2	1:E:66:ARG:NH1	2.29	0.47
1:H:192:ASP:O	1:H:213:ALA:HA	2.15	0.46
1:G:139:TRP:CD2	1:G:161:LEU:HG	2.50	0.46
1:A:3:VAL:HG21	1:A:84:PRO:HB3	1.97	0.46
1:C:3:VAL:HG13	1:C:4:ILE:HD12	1.96	0.46
1:E:63:GYC:HA31	1:E:63:GYC:N	2.31	0.46
1:G:156:ASP:OD2	2:G:302:HOH:O	2.21	0.46
1:E:59:THR:HG22	1:E:63:GYC:CE2	2.46	0.46
1:G:60:VAL:HG13	1:G:89:TRP:HH2	1.81	0.46
1:G:9:LYS:HE2	1:G:112:ASP:OD2	2.16	0.45
1:H:13:ARG:NH2	1:H:15:GLU:OE1	2.40	0.45
1:D:146:LEU:HB3	1:D:153:LEU:HD11	1.99	0.45
1:D:174:LYS:HG3	2:D:309:HOH:O	2.16	0.45
1:E:191:VAL:HA	1:E:215:ALA:HA	1.98	0.45
1:F:198:LYS:HD3	1:F:210:HIS:CD2	2.52	0.45
1:H:63:GYC:HB2	1:H:66:ARG:NH2	2.30	0.45
1:C:102:ASN:HD21	1:E:100:ILE:HG22	1.81	0.45
1:D:9:LYS:O	1:D:113:CYS:HA	2.16	0.45
1:D:149:ARG:O	1:D:152:VAL:HG22	2.16	0.45
1:A:146:LEU:HA	1:A:154:LYS:O	2.16	0.45
1:D:25:ILE:HG12	1:D:44:VAL:HG22	1.98	0.45
1:G:196:GLU:OE1	1:G:198:LYS:HE2	2.17	0.45
1:F:89:TRP:CZ3	1:F:91:ARG:HB3	2.52	0.45
1:G:100:ILE:O	1:G:122:GLY:HA2	2.17	0.44
1:A:153:LEU:HB3	1:A:177:TYR:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:81:GLN:HG2	1:E:183:VAL:HG12	1.99	0.44
1:A:43:LYS:HD2	1:A:206:ASN:OD1	2.16	0.44
1:G:158:ASN:N	1:G:158:ASN:OD1	2.50	0.44
1:B:146:LEU:HB3	1:B:153:LEU:HD11	2.00	0.44
1:D:170:ARG:NH2	1:H:149:ARG:HH12	2.16	0.44
1:E:145:LYS:HB2	1:E:145:LYS:HE3	1.54	0.44
1:G:90:GLU:CG	1:G:176:THR:HB	2.47	0.44
1:H:144:GLU:HB2	1:H:157:VAL:CG2	2.47	0.44
1:H:144:GLU:CB	1:H:157:VAL:CG2	2.96	0.44
1:C:33:PRO:O	1:C:80:LYS:HE2	2.15	0.44
1:E:63:GYC:HB12	1:E:195:ILE:HD12	1.99	0.44
1:D:146:LEU:HA	1:D:154:LYS:O	2.17	0.43
1:E:93:MET:HB2	1:E:101:CYS:HB2	2.00	0.43
1:F:146:LEU:HD13	1:F:153:LEU:HD21	2.00	0.43
1:E:140:GLU:OE2	1:E:168:HIS:NE2	2.34	0.43
1:G:139:TRP:CE2	1:G:161:LEU:CG	3.01	0.43
1:H:93:MET:HB2	1:H:101:CYS:HB2	2.00	0.43
1:H:219:LEU:HA	1:H:220:PRO:HD3	1.88	0.43
1:A:137:VAL:HG22	1:A:164:GLU:HG2	2.00	0.43
1:H:97:ASP:OD1	1:H:169:TYR:OH	2.31	0.43
1:F:77:ASP:OD1	1:F:80:LYS:HD2	2.19	0.43
1:H:36:GLY:O	1:H:212:HIS:HA	2.19	0.43
1:B:171:CYS:HG	1:B:173:PHE:HE1	1.64	0.43
1:A:71:TYR:HA	1:A:72:PRO:HD3	1.86	0.43
1:E:94[A]:ASN:OD1	1:E:94[A]:ASN:N	2.52	0.43
1:H:21:HIS:HA	1:H:22:PRO:HD3	1.92	0.43
1:C:138:LYS:HB2	1:C:138:LYS:HE3	1.90	0.43
1:H:200:HIS:HA	1:H:206:ASN:O	2.19	0.43
1:D:93:MET:HB2	1:D:101:CYS:HB2	2.00	0.43
1:D:192:ASP:HB3	1:H:219:LEU:CD1	2.48	0.43
1:E:144:GLU:CA	1:E:157:VAL:HG12	2.49	0.43
1:G:63:GYC:HB11	1:G:195:ILE:HD12	2.00	0.43
1:B:158:ASN:O	1:F:145:LYS:HD2	2.19	0.42
1:A:81:GLN:HB3	1:A:183:VAL:HG13	2.01	0.42
1:A:192:ASP:O	1:A:213:ALA:HA	2.19	0.42
1:C:192:ASP:O	1:C:213:ALA:HA	2.18	0.42
1:H:81:GLN:HB3	1:H:183:VAL:HG11	2.01	0.42
1:D:39:SER:HB2	1:D:210:HIS:CE1	2.54	0.42
1:F:192:ASP:O	1:F:213:ALA:HA	2.19	0.42
1:G:145:LYS:HB2	1:G:145:LYS:HE2	1.83	0.42
1:B:145:LYS:NZ	1:B:156:ASP:OD1	2.36	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:157:VAL:HB	1:F:173:PHE:CD2	2.54	0.42
1:B:85:GLU:CD	1:B:181:LYS:HD2	2.40	0.42
1:F:145:LYS:HE2	1:F:145:LYS:HB2	1.78	0.42
1:H:183:VAL:HG12	1:H:184:GLN:H	1.85	0.42
1:A:77:ASP:O	1:A:81:GLN:HB2	2.19	0.42
1:E:219:LEU:HA	1:E:220:PRO:HD3	1.78	0.42
1:G:142:ALA:HB2	1:G:159:THR:HG21	2.01	0.42
1:C:52:PHE:HE1	1:C:57:LEU:HD11	1.84	0.42
1:D:96:GLU:OE1	1:H:149:ARG:NH2	2.47	0.42
1:E:198:LYS:HG3	1:E:210:HIS:CD2	2.55	0.42
1:F:27:GLY:HA3	1:F:42:LEU:HD23	2.01	0.42
1:A:159:THR:OG1	1:A:160:ALA:N	2.53	0.42
1:D:85:GLU:HB2	1:D:181:LYS:HD3	2.02	0.42
1:A:63:GYC:HB2	1:A:66:ARG:HH22	1.83	0.41
1:E:63:GYC:CB1	1:E:195:ILE:HD12	2.50	0.41
1:H:157:VAL:HB	1:H:173:PHE:HB2	2.02	0.41
1:B:100:ILE:O	1:B:122:GLY:HA2	2.19	0.41
1:B:153:LEU:HB3	1:B:177:TYR:HB2	2.02	0.41
1:C:32:LYS:HB3	1:C:35:GLU:HG3	2.02	0.41
1:F:53:ALA:O	1:F:56:ILE:HG12	2.20	0.41
1:F:89:TRP:HZ3	1:F:91:ARG:HB3	1.86	0.41
1:G:60:VAL:HG13	1:G:89:TRP:CH2	2.55	0.41
1:B:33:PRO:HA	1:B:68:PHE:HA	2.01	0.41
1:C:94:ASN:OD1	1:C:100:ILE:HD12	2.20	0.41
1:E:36:GLY:O	1:E:212:HIS:HA	2.21	0.41
1:A:144:GLU:HB2	1:A:157:VAL:HG22	2.03	0.41
1:C:3:VAL:CG1	1:C:4:ILE:HD12	2.50	0.41
1:H:146:LEU:N	1:H:146:LEU:HD12	2.36	0.41
1:B:139:TRP:CZ3	1:B:159:THR:HG23	2.56	0.41
1:A:12:LEU:HB2	1:A:116:TYR:HB2	2.03	0.41
1:A:183:VAL:HG12	1:A:184:GLN:N	2.36	0.41
1:E:146:LEU:HA	1:E:154:LYS:O	2.21	0.41
1:A:110:ASP:OD2	1:F:45:LYS:NZ	2.53	0.41
1:D:203:ASP:OD2	1:D:205:SER:OG	2.39	0.41
1:F:33:PRO:HA	1:F:68:PHE:HA	2.02	0.41
1:H:128:ASN:HA	1:H:133:GLN:HE21	1.86	0.41
1:F:219:LEU:HA	1:F:220:PRO:HD3	1.92	0.40
1:A:42:LEU:O	1:A:43:LYS:HD3	2.20	0.40
1:B:157:VAL:HB	1:B:173:PHE:CB	2.48	0.40
1:C:34:PHE:CE1	1:C:80:LYS:HE2	2.57	0.40
1:B:78:TYR:HB2	1:B:186:PRO:HD3	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:145:LYS:HD2	1:G:158:ASN:O	2.21	0.40
1:D:143:THR:CB	1:H:143:THR:HG21	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	213/255 (84%)	209 (98%)	4 (2%)	0	100	100
1	B	213/255 (84%)	209 (98%)	4 (2%)	0	100	100
1	C	216/255 (85%)	211 (98%)	4 (2%)	1 (0%)	29	43
1	D	215/255 (84%)	211 (98%)	4 (2%)	0	100	100
1	E	215/255 (84%)	211 (98%)	4 (2%)	0	100	100
1	F	214/255 (84%)	209 (98%)	5 (2%)	0	100	100
1	G	213/255 (84%)	206 (97%)	7 (3%)	0	100	100
1	H	213/255 (84%)	210 (99%)	3 (1%)	0	100	100
All	All	1712/2040 (84%)	1676 (98%)	35 (2%)	1 (0%)	51	69

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	33	PRO

5.3.2 Protein sidechains [i](#)

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	181/216 (84%)	173 (96%)	8 (4%)	28	44
1	B	182/216 (84%)	175 (96%)	7 (4%)	33	50
1	C	179/216 (83%)	172 (96%)	7 (4%)	32	48
1	D	180/216 (83%)	176 (98%)	4 (2%)	52	70
1	E	181/216 (84%)	175 (97%)	6 (3%)	38	55
1	F	185/216 (86%)	176 (95%)	9 (5%)	25	39
1	G	180/216 (83%)	176 (98%)	4 (2%)	52	70
1	H	177/216 (82%)	172 (97%)	5 (3%)	43	61
All	All	1445/1728 (84%)	1395 (96%)	50 (4%)	36	53

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASP
1	A	15	GLU
1	A	30	LEU
1	A	38	GLN
1	A	154	LYS
1	A	180	LYS
1	A	181	LYS
1	A	202	LYS
1	B	66	ARG
1	B	132	MET
1	B	158	ASN
1	B	162	SER
1	B	180	LYS
1	B	181	LYS
1	B	211	GLU
1	C	30	LEU
1	C	42	LEU
1	C	43	LYS
1	C	66	ARG
1	C	94	ASN
1	C	161	LEU
1	C	212	HIS
1	D	6	PRO
1	D	144	GLU

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Mol	Chain	Res	Type
1	D	158	ASN
1	D	181	LYS
1	E	43	LYS
1	E	66	ARG
1	E	74	ASN
1	E	134	LYS
1	E	145	LYS
1	E	181	LYS
1	F	32	LYS
1	F	38	GLN
1	F	39	SER
1	F	66	ARG
1	F	91	ARG
1	F	94	ASN
1	F	117	GLU
1	F	162	SER
1	F	164	GLU
1	G	38	GLN
1	G	39	SER
1	G	43	LYS
1	G	159	THR
1	H	28	VAL
1	H	30	LEU
1	H	66	ARG
1	H	132	MET
1	H	211	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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	GYC	E	63	1	22,22,23	1.94	7 (31%)	26,30,32	2.35	7 (26%)
1	GYC	B	63	1	22,22,23	1.15	1 (4%)	26,30,32	2.70	7 (26%)
1	GYC	G	63	1	22,22,23	1.18	1 (4%)	26,30,32	2.62	8 (30%)
1	GYC	F	63	1	22,22,23	1.14	1 (4%)	26,30,32	2.71	8 (30%)
1	GYC	C	63	1	22,22,23	1.23	1 (4%)	26,30,32	2.49	7 (26%)
1	GYC	D	63	1	22,22,23	1.13	1 (4%)	26,30,32	2.62	7 (26%)
1	GYC	H	63	1	22,22,23	1.11	1 (4%)	26,30,32	2.63	8 (30%)
1	GYC	A	63	1	22,22,23	1.81	8 (36%)	26,30,32	2.35	5 (19%)

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	GYC	E	63	1	-	4/9/29/30	0/2/2/2
1	GYC	B	63	1	-	4/9/29/30	0/2/2/2
1	GYC	G	63	1	-	3/9/29/30	0/2/2/2
1	GYC	F	63	1	-	4/9/29/30	0/2/2/2
1	GYC	C	63	1	-	3/9/29/30	0/2/2/2
1	GYC	D	63	1	-	3/9/29/30	0/2/2/2
1	GYC	H	63	1	-	4/9/29/30	0/2/2/2
1	GYC	A	63	1	-	3/9/29/30	0/2/2/2

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	63	GYC	CB2-CA2	4.93	1.39	1.35
1	G	63	GYC	CB2-CA2	4.65	1.39	1.35
1	B	63	GYC	CB2-CA2	4.53	1.38	1.35
1	F	63	GYC	CB2-CA2	4.47	1.38	1.35
1	H	63	GYC	CB2-CA2	4.36	1.38	1.35
1	D	63	GYC	CB2-CA2	4.27	1.38	1.35
1	E	63	GYC	CB2-CA2	4.20	1.38	1.35
1	A	63	GYC	CA2-C2	-3.86	1.44	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	63	GYC	C2-N3	-3.46	1.31	1.39
1	A	63	GYC	C2-N3	-3.24	1.32	1.39
1	E	63	GYC	C1-N2	-2.91	1.27	1.32
1	E	63	GYC	CA2-C2	-2.86	1.45	1.48
1	E	63	GYC	CB1-CA1	-2.72	1.50	1.53
1	E	63	GYC	C1-N3	-2.52	1.32	1.37
1	A	63	GYC	C1-N2	-2.46	1.28	1.32
1	E	63	GYC	CE2-CD2	-2.24	1.34	1.38
1	A	63	GYC	CA2-N2	-2.21	1.33	1.38
1	A	63	GYC	CE1-CD1	-2.16	1.34	1.38
1	A	63	GYC	CE2-CD2	-2.14	1.34	1.38
1	A	63	GYC	CE1-CZ	-2.08	1.35	1.38
1	A	63	GYC	CD2-CG2	-2.00	1.35	1.39

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	63	GYC	O2-C2-CA2	-8.11	126.41	130.96
1	F	63	GYC	O2-C2-CA2	-7.65	126.66	130.96
1	D	63	GYC	O2-C2-CA2	-7.51	126.74	130.96
1	H	63	GYC	O2-C2-CA2	-7.40	126.80	130.96
1	G	63	GYC	O2-C2-CA2	-6.85	127.11	130.96
1	C	63	GYC	O2-C2-CA2	-6.68	127.21	130.96
1	E	63	GYC	C2-CA2-N2	-6.41	104.45	108.93
1	A	63	GYC	O-C-CA3	-5.70	109.17	126.39
1	E	63	GYC	CA2-C2-N3	5.51	105.97	103.37
1	G	63	GYC	C2-CA2-N2	-5.39	105.16	108.93
1	H	63	GYC	CA2-C2-N3	5.30	105.88	103.37
1	C	63	GYC	C2-CA2-N2	-5.29	105.23	108.93
1	F	63	GYC	C2-CA2-N2	-5.29	105.23	108.93
1	G	63	GYC	CA2-N2-C1	5.25	109.64	105.77
1	A	63	GYC	CA2-N2-C1	5.19	109.60	105.77
1	B	63	GYC	C2-CA2-N2	-5.19	105.30	108.93
1	A	63	GYC	O2-C2-CA2	-5.16	128.06	130.96
1	F	63	GYC	CA2-C2-N3	5.13	105.80	103.37
1	E	63	GYC	CA2-N2-C1	5.06	109.50	105.77
1	B	63	GYC	CA2-C2-N3	5.02	105.75	103.37
1	G	63	GYC	CA2-C2-N3	4.99	105.73	103.37
1	H	63	GYC	C2-CA2-N2	-4.95	105.47	108.93
1	B	63	GYC	CA2-N2-C1	4.94	109.41	105.77
1	F	63	GYC	CG2-CB2-CA2	-4.92	123.91	129.94
1	D	63	GYC	C2-CA2-N2	-4.92	105.49	108.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	63	GYC	CG2-CB2-CA2	-4.92	123.92	129.94
1	F	63	GYC	CA2-N2-C1	4.90	109.38	105.77
1	D	63	GYC	CA2-N2-C1	4.87	109.36	105.77
1	A	63	GYC	CA2-C2-N3	4.86	105.67	103.37
1	D	63	GYC	CA2-C2-N3	4.83	105.66	103.37
1	C	63	GYC	CA2-N2-C1	4.74	109.26	105.77
1	C	63	GYC	CA2-C2-N3	4.71	105.60	103.37
1	H	63	GYC	CA2-N2-C1	4.40	109.02	105.77
1	G	63	GYC	CG2-CB2-CA2	-4.35	124.61	129.94
1	H	63	GYC	CG2-CB2-CA2	-4.29	124.69	129.94
1	A	63	GYC	C2-CA2-N2	-4.20	105.99	108.93
1	E	63	GYC	O2-C2-CA2	-4.01	128.71	130.96
1	B	63	GYC	CG2-CB2-CA2	-3.87	125.21	129.94
1	H	63	GYC	CA1-CB1-SG1	-3.39	107.14	114.44
1	D	63	GYC	CB2-CA2-C2	3.21	126.11	122.28
1	C	63	GYC	CG2-CB2-CA2	-3.18	126.05	129.94
1	G	63	GYC	CB2-CA2-C2	2.96	125.81	122.28
1	E	63	GYC	O-C-CA3	-2.89	117.67	126.39
1	E	63	GYC	CA1-CB1-SG1	-2.86	108.27	114.44
1	F	63	GYC	CB2-CA2-C2	2.78	125.60	122.28
1	C	63	GYC	CA3-N3-C1	-2.61	124.03	127.16
1	C	63	GYC	O-C-CA3	-2.57	118.64	126.39
1	F	63	GYC	CA1-CB1-SG1	-2.55	108.95	114.44
1	B	63	GYC	CA1-CB1-SG1	-2.50	109.05	114.44
1	G	63	GYC	O-C-CA3	-2.41	119.12	126.39
1	H	63	GYC	CB2-CA2-C2	2.37	125.11	122.28
1	H	63	GYC	O-C-CA3	-2.31	119.42	126.39
1	D	63	GYC	CA1-CB1-SG1	-2.29	109.50	114.44
1	B	63	GYC	O-C-CA3	-2.26	119.55	126.39
1	G	63	GYC	CA1-CB1-SG1	-2.13	109.84	114.44
1	E	63	GYC	CB2-CA2-C2	2.05	124.72	122.28
1	F	63	GYC	O-C-CA3	-2.02	120.29	126.39

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	63	GYC	C2-CA2-CB2-CG2
1	B	63	GYC	C-CA3-N3-C2
1	B	63	GYC	N2-CA2-CB2-CG2
1	B	63	GYC	C2-CA2-CB2-CG2
1	C	63	GYC	C-CA3-N3-C2

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Mol	Chain	Res	Type	Atoms
1	C	63	GYC	N2-CA2-CB2-CG2
1	C	63	GYC	C2-CA2-CB2-CG2
1	D	63	GYC	C-CA3-N3-C2
1	D	63	GYC	C2-CA2-CB2-CG2
1	E	63	GYC	C2-CA2-CB2-CG2
1	F	63	GYC	C-CA3-N3-C2
1	F	63	GYC	N2-CA2-CB2-CG2
1	F	63	GYC	C2-CA2-CB2-CG2
1	G	63	GYC	C-CA3-N3-C2
1	G	63	GYC	N2-CA2-CB2-CG2
1	G	63	GYC	C2-CA2-CB2-CG2
1	H	63	GYC	C-CA3-N3-C2
1	H	63	GYC	N2-CA2-CB2-CG2
1	H	63	GYC	C2-CA2-CB2-CG2
1	A	63	GYC	N2-CA2-CB2-CG2
1	D	63	GYC	N2-CA2-CB2-CG2
1	E	63	GYC	N2-CA2-CB2-CG2
1	A	63	GYC	C-CA3-N3-C2
1	E	63	GYC	C1-CA1-CB1-SG1
1	B	63	GYC	C-CA3-N3-C1
1	F	63	GYC	C-CA3-N3-C1
1	H	63	GYC	C-CA3-N3-C1
1	E	63	GYC	C-CA3-N3-C2

There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	E	63	GYC	5	0
1	G	63	GYC	1	0
1	F	63	GYC	1	0
1	H	63	GYC	1	0
1	A	63	GYC	2	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	215/255 (84%)	0.22	6 (2%)	53	49	18, 34, 60, 70	0
1	B	215/255 (84%)	0.22	8 (3%)	41	38	20, 33, 61, 79	0
1	C	220/255 (86%)	0.48	11 (5%)	28	25	22, 42, 67, 91	0
1	D	219/255 (85%)	0.22	6 (2%)	54	50	19, 36, 60, 73	0
1	E	218/255 (85%)	0.40	11 (5%)	28	25	23, 40, 59, 74	0
1	F	218/255 (85%)	0.25	6 (2%)	53	49	18, 34, 60, 77	0
1	G	215/255 (84%)	0.26	6 (2%)	53	49	21, 37, 63, 91	0
1	H	215/255 (84%)	0.26	7 (3%)	46	43	21, 36, 61, 77	0
All	All	1735/2040 (85%)	0.29	61 (3%)	44	40	18, 37, 62, 91	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	150	ASP	4.7
1	C	224	MET	4.4
1	F	3	VAL	4.3
1	C	3	VAL	4.2
1	G	3	VAL	4.1
1	H	111	GLY	3.8
1	H	220	PRO	3.7
1	B	165	GLY	3.2
1	B	110	ASP	3.2
1	D	201	ASP	3.2
1	F	220	PRO	3.1
1	E	201	ASP	3.1
1	F	151	GLY	3.0
1	E	111	GLY	3.0
1	G	166	GLY	3.0
1	E	3	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	F	112	ASP	2.9
1	F	150	ASP	2.9
1	B	166	GLY	2.8
1	C	7	ASP	2.8
1	E	164	GLU	2.7
1	C	21	HIS	2.7
1	A	158	ASN	2.6
1	H	182	VAL	2.6
1	G	110	ASP	2.6
1	E	48	GLY	2.6
1	C	227	LEU	2.5
1	A	166	GLY	2.5
1	D	3	VAL	2.5
1	D	128	ASN	2.5
1	E	165	GLY	2.5
1	F	110	ASP	2.5
1	D	212	HIS	2.5
1	E	203	ASP	2.4
1	B	111	GLY	2.4
1	H	158	ASN	2.3
1	E	112	ASP	2.3
1	D	127	ALA	2.3
1	C	220	PRO	2.2
1	E	202	LYS	2.2
1	B	220	PRO	2.2
1	B	5	LYS	2.2
1	C	10	ILE	2.2
1	A	111	GLY	2.2
1	H	3	VAL	2.1
1	C	6	PRO	2.1
1	C	50	LEU	2.1
1	E	110	ASP	2.1
1	B	4	ILE	2.1
1	A	110	ASP	2.1
1	G	7	ASP	2.1
1	D	150	ASP	2.1
1	B	150	ASP	2.1
1	A	220	PRO	2.1
1	H	136	THR	2.1
1	H	213	ALA	2.1
1	C	111	GLY	2.1
1	G	182	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	196	GLU	2.0
1	G	112	ASP	2.0
1	A	5	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	GYC	E	63	21/22	0.72	0.29	27,41,56,63	0
1	GYC	G	63	21/22	0.90	0.22	24,38,57,58	0
1	GYC	C	63	21/22	0.91	0.16	29,32,39,52	0
1	GYC	A	63	21/22	0.91	0.15	23,42,48,70	0
1	GYC	B	63	21/22	0.91	0.21	15,26,37,40	0
1	GYC	F	63	21/22	0.92	0.16	19,27,34,43	0
1	GYC	H	63	21/22	0.92	0.18	7,25,42,50	0
1	GYC	D	63	21/22	0.93	0.15	22,29,42,63	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.