



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:13 PM UTC

PDB ID : 7Z7P / pdb_00007z7p
Title : Structure of the fluorescent protein NeonCyan0.95 at pH 5.6
Authors : Depernet, H.; Dupuy, J.; Royant, A.
Deposited on : 2022-03-16
Resolution : 1.95 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

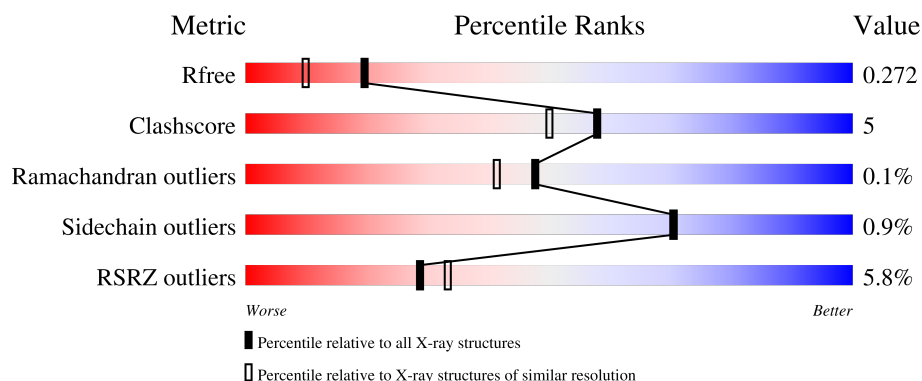
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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}	180053	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	234	4% 83% 13% .
1	B	234	5% 82% 13% 5%
1	C	234	4% 86% 10% .
1	D	234	6% 90% 7% .
1	E	234	8% 84% 9% 6%

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Mol	Chain	Length	Quality of chain
1	F	234	6% 85% 9% 6%
1	G	234	7% 85% 8% 6%
1	H	234	3% 80% 13% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14907 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 NeonCyan0.95.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	6	0
			1842	1175	308	348	11			
1	B	223	Total	C	N	O	S	0	6	0
			1813	1153	304	346	10			
1	C	227	Total	C	N	O	S	0	6	0
			1849	1177	309	353	10			
1	D	227	Total	C	N	O	S	0	6	0
			1841	1170	308	353	10			
1	E	220	Total	C	N	O	S	0	1	0
			1749	1116	291	334	8			
1	F	221	Total	C	N	O	S	0	4	0
			1770	1132	295	334	9			
1	G	219	Total	C	N	O	S	0	3	0
			1752	1119	292	332	9			
1	H	220	Total	C	N	O	S	0	2	0
			1755	1121	294	332	8			

There are 392 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP B1PNC0
A	2	VAL	-	expression tag	UNP B1PNC0
A	3	SER	-	expression tag	UNP B1PNC0
A	4	LYS	-	expression tag	UNP B1PNC0
A	5	GLY	-	expression tag	UNP B1PNC0
A	6	GLU	-	expression tag	UNP B1PNC0
A	7	GLU	-	expression tag	UNP B1PNC0
A	8	ASP	-	expression tag	UNP B1PNC0
A	9	ASN	-	expression tag	UNP B1PNC0
A	10	MET	-	expression tag	UNP B1PNC0
A	11	ALA	-	expression tag	UNP B1PNC0
A	25	ILE	PHE	engineered mutation	UNP B1PNC0
A	35	GLN	ARG	engineered mutation	UNP B1PNC0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	42	GLU	ASP	engineered mutation	UNP B1PNC0
A	55	ASP	ALA	engineered mutation	UNP B1PNC0
A	66	HIS	GLN	engineered mutation	UNP B1PNC0
A	67	VAL	ILE	engineered mutation	UNP B1PNC0
A	69	IO8	GLY	chromophore	UNP B1PNC0
A	?	-	TYR	chromophore	UNP B1PNC0
A	?	-	GLY	chromophore	UNP B1PNC0
A	77	TYR	PHE	engineered mutation	UNP B1PNC0
A	89	VAL	LYS	engineered mutation	UNP B1PNC0
A	110	VAL	SER	engineered mutation	UNP B1PNC0
A	125	ALA	PHE	engineered mutation	UNP B1PNC0
A	128	LYS	ILE	engineered mutation	UNP B1PNC0
A	146	VAL	ALA	engineered mutation	UNP B1PNC0
A	148	SER	TRP	engineered mutation	UNP B1PNC0
A	150	TRP	VAL	engineered mutation	UNP B1PNC0
A	151	SER	THR	engineered mutation	UNP B1PNC0
A	153	LYS	MET	engineered mutation	UNP B1PNC0
A	154	THR	LEU	engineered mutation	UNP B1PNC0
A	166	LYS	ASP	engineered mutation	UNP B1PNC0
A	168	SER	THR	engineered mutation	UNP B1PNC0
A	173	ASN	SER	engineered mutation	UNP B1PNC0
A	175	GLU	LYS	engineered mutation	UNP B1PNC0
A	178	ARG	GLN	engineered mutation	UNP B1PNC0
A	181	ALA	VAL	engineered mutation	UNP B1PNC0
A	184	THR	ASN	engineered mutation	UNP B1PNC0
A	195	TYR	ILE	engineered mutation	UNP B1PNC0
A	201	VAL	MET	engineered mutation	UNP B1PNC0
A	202	TYR	PHE	engineered mutation	UNP B1PNC0
A	213	GLU	LYS	engineered mutation	UNP B1PNC0
A	230	GLY	-	expression tag	UNP B1PNC0
A	231	MET	-	expression tag	UNP B1PNC0
A	232	ASP	-	expression tag	UNP B1PNC0
A	233	GLU	-	expression tag	UNP B1PNC0
A	234	LEU	-	expression tag	UNP B1PNC0
A	235	TYR	-	expression tag	UNP B1PNC0
A	236	LYS	-	expression tag	UNP B1PNC0
B	1	MET	-	initiating methionine	UNP B1PNC0
B	2	VAL	-	expression tag	UNP B1PNC0
B	3	SER	-	expression tag	UNP B1PNC0
B	4	LYS	-	expression tag	UNP B1PNC0
B	5	GLY	-	expression tag	UNP B1PNC0
B	6	GLU	-	expression tag	UNP B1PNC0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	7	GLU	-	expression tag	UNP B1PNC0
B	8	ASP	-	expression tag	UNP B1PNC0
B	9	ASN	-	expression tag	UNP B1PNC0
B	10	MET	-	expression tag	UNP B1PNC0
B	11	ALA	-	expression tag	UNP B1PNC0
B	25	ILE	PHE	engineered mutation	UNP B1PNC0
B	35	GLN	ARG	engineered mutation	UNP B1PNC0
B	42	GLU	ASP	engineered mutation	UNP B1PNC0
B	55	ASP	ALA	engineered mutation	UNP B1PNC0
B	66	HIS	GLN	engineered mutation	UNP B1PNC0
B	67	VAL	ILE	engineered mutation	UNP B1PNC0
B	69	IO8	GLY	chromophore	UNP B1PNC0
B	?	-	TYR	chromophore	UNP B1PNC0
B	?	-	GLY	chromophore	UNP B1PNC0
B	77	TYR	PHE	engineered mutation	UNP B1PNC0
B	89	VAL	LYS	engineered mutation	UNP B1PNC0
B	110	VAL	SER	engineered mutation	UNP B1PNC0
B	125	ALA	PHE	engineered mutation	UNP B1PNC0
B	128	LYS	ILE	engineered mutation	UNP B1PNC0
B	146	VAL	ALA	engineered mutation	UNP B1PNC0
B	148	SER	TRP	engineered mutation	UNP B1PNC0
B	150	TRP	VAL	engineered mutation	UNP B1PNC0
B	151	SER	THR	engineered mutation	UNP B1PNC0
B	153	LYS	MET	engineered mutation	UNP B1PNC0
B	154	THR	LEU	engineered mutation	UNP B1PNC0
B	166	LYS	ASP	engineered mutation	UNP B1PNC0
B	168	SER	THR	engineered mutation	UNP B1PNC0
B	173	ASN	SER	engineered mutation	UNP B1PNC0
B	175	GLU	LYS	engineered mutation	UNP B1PNC0
B	178	ARG	GLN	engineered mutation	UNP B1PNC0
B	181	ALA	VAL	engineered mutation	UNP B1PNC0
B	184	THR	ASN	engineered mutation	UNP B1PNC0
B	195	TYR	ILE	engineered mutation	UNP B1PNC0
B	201	VAL	MET	engineered mutation	UNP B1PNC0
B	202	TYR	PHE	engineered mutation	UNP B1PNC0
B	213	GLU	LYS	engineered mutation	UNP B1PNC0
B	230	GLY	-	expression tag	UNP B1PNC0
B	231	MET	-	expression tag	UNP B1PNC0
B	232	ASP	-	expression tag	UNP B1PNC0
B	233	GLU	-	expression tag	UNP B1PNC0
B	234	LEU	-	expression tag	UNP B1PNC0
B	235	TYR	-	expression tag	UNP B1PNC0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	236	LYS	-	expression tag	UNP B1PNC0
C	1	MET	-	initiating methionine	UNP B1PNC0
C	2	VAL	-	expression tag	UNP B1PNC0
C	3	SER	-	expression tag	UNP B1PNC0
C	4	LYS	-	expression tag	UNP B1PNC0
C	5	GLY	-	expression tag	UNP B1PNC0
C	6	GLU	-	expression tag	UNP B1PNC0
C	7	GLU	-	expression tag	UNP B1PNC0
C	8	ASP	-	expression tag	UNP B1PNC0
C	9	ASN	-	expression tag	UNP B1PNC0
C	10	MET	-	expression tag	UNP B1PNC0
C	11	ALA	-	expression tag	UNP B1PNC0
C	25	ILE	PHE	engineered mutation	UNP B1PNC0
C	35	GLN	ARG	engineered mutation	UNP B1PNC0
C	42	GLU	ASP	engineered mutation	UNP B1PNC0
C	55	ASP	ALA	engineered mutation	UNP B1PNC0
C	66	HIS	GLN	engineered mutation	UNP B1PNC0
C	67	VAL	ILE	engineered mutation	UNP B1PNC0
C	69	IO8	GLY	chromophore	UNP B1PNC0
C	?	-	TYR	chromophore	UNP B1PNC0
C	?	-	GLY	chromophore	UNP B1PNC0
C	77	TYR	PHE	engineered mutation	UNP B1PNC0
C	89	VAL	LYS	engineered mutation	UNP B1PNC0
C	110	VAL	SER	engineered mutation	UNP B1PNC0
C	125	ALA	PHE	engineered mutation	UNP B1PNC0
C	128	LYS	ILE	engineered mutation	UNP B1PNC0
C	146	VAL	ALA	engineered mutation	UNP B1PNC0
C	148	SER	TRP	engineered mutation	UNP B1PNC0
C	150	TRP	VAL	engineered mutation	UNP B1PNC0
C	151	SER	THR	engineered mutation	UNP B1PNC0
C	153	LYS	MET	engineered mutation	UNP B1PNC0
C	154	THR	LEU	engineered mutation	UNP B1PNC0
C	166	LYS	ASP	engineered mutation	UNP B1PNC0
C	168	SER	THR	engineered mutation	UNP B1PNC0
C	173	ASN	SER	engineered mutation	UNP B1PNC0
C	175	GLU	LYS	engineered mutation	UNP B1PNC0
C	178	ARG	GLN	engineered mutation	UNP B1PNC0
C	181	ALA	VAL	engineered mutation	UNP B1PNC0
C	184	THR	ASN	engineered mutation	UNP B1PNC0
C	195	TYR	ILE	engineered mutation	UNP B1PNC0
C	201	VAL	MET	engineered mutation	UNP B1PNC0
C	202	TYR	PHE	engineered mutation	UNP B1PNC0

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Chain	Residue	Modelled	Actual	Comment	Reference
C	213	GLU	LYS	engineered mutation	UNP B1PNC0
C	230	GLY	-	expression tag	UNP B1PNC0
C	231	MET	-	expression tag	UNP B1PNC0
C	232	ASP	-	expression tag	UNP B1PNC0
C	233	GLU	-	expression tag	UNP B1PNC0
C	234	LEU	-	expression tag	UNP B1PNC0
C	235	TYR	-	expression tag	UNP B1PNC0
C	236	LYS	-	expression tag	UNP B1PNC0
D	1	MET	-	initiating methionine	UNP B1PNC0
D	2	VAL	-	expression tag	UNP B1PNC0
D	3	SER	-	expression tag	UNP B1PNC0
D	4	LYS	-	expression tag	UNP B1PNC0
D	5	GLY	-	expression tag	UNP B1PNC0
D	6	GLU	-	expression tag	UNP B1PNC0
D	7	GLU	-	expression tag	UNP B1PNC0
D	8	ASP	-	expression tag	UNP B1PNC0
D	9	ASN	-	expression tag	UNP B1PNC0
D	10	MET	-	expression tag	UNP B1PNC0
D	11	ALA	-	expression tag	UNP B1PNC0
D	25	ILE	PHE	engineered mutation	UNP B1PNC0
D	35	GLN	ARG	engineered mutation	UNP B1PNC0
D	42	GLU	ASP	engineered mutation	UNP B1PNC0
D	55	ASP	ALA	engineered mutation	UNP B1PNC0
D	66	HIS	GLN	engineered mutation	UNP B1PNC0
D	67	VAL	ILE	engineered mutation	UNP B1PNC0
D	69	IO8	GLY	chromophore	UNP B1PNC0
D	?	-	TYR	chromophore	UNP B1PNC0
D	?	-	GLY	chromophore	UNP B1PNC0
D	77	TYR	PHE	engineered mutation	UNP B1PNC0
D	89	VAL	LYS	engineered mutation	UNP B1PNC0
D	110	VAL	SER	engineered mutation	UNP B1PNC0
D	125	ALA	PHE	engineered mutation	UNP B1PNC0
D	128	LYS	ILE	engineered mutation	UNP B1PNC0
D	146	VAL	ALA	engineered mutation	UNP B1PNC0
D	148	SER	TRP	engineered mutation	UNP B1PNC0
D	150	TRP	VAL	engineered mutation	UNP B1PNC0
D	151	SER	THR	engineered mutation	UNP B1PNC0
D	153	LYS	MET	engineered mutation	UNP B1PNC0
D	154	THR	LEU	engineered mutation	UNP B1PNC0
D	166	LYS	ASP	engineered mutation	UNP B1PNC0
D	168	SER	THR	engineered mutation	UNP B1PNC0
D	173	ASN	SER	engineered mutation	UNP B1PNC0

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Chain	Residue	Modelled	Actual	Comment	Reference
D	175	GLU	LYS	engineered mutation	UNP B1PNC0
D	178	ARG	GLN	engineered mutation	UNP B1PNC0
D	181	ALA	VAL	engineered mutation	UNP B1PNC0
D	184	THR	ASN	engineered mutation	UNP B1PNC0
D	195	TYR	ILE	engineered mutation	UNP B1PNC0
D	201	VAL	MET	engineered mutation	UNP B1PNC0
D	202	TYR	PHE	engineered mutation	UNP B1PNC0
D	213	GLU	LYS	engineered mutation	UNP B1PNC0
D	230	GLY	-	expression tag	UNP B1PNC0
D	231	MET	-	expression tag	UNP B1PNC0
D	232	ASP	-	expression tag	UNP B1PNC0
D	233	GLU	-	expression tag	UNP B1PNC0
D	234	LEU	-	expression tag	UNP B1PNC0
D	235	TYR	-	expression tag	UNP B1PNC0
D	236	LYS	-	expression tag	UNP B1PNC0
E	1	MET	-	initiating methionine	UNP B1PNC0
E	2	VAL	-	expression tag	UNP B1PNC0
E	3	SER	-	expression tag	UNP B1PNC0
E	4	LYS	-	expression tag	UNP B1PNC0
E	5	GLY	-	expression tag	UNP B1PNC0
E	6	GLU	-	expression tag	UNP B1PNC0
E	7	GLU	-	expression tag	UNP B1PNC0
E	8	ASP	-	expression tag	UNP B1PNC0
E	9	ASN	-	expression tag	UNP B1PNC0
E	10	MET	-	expression tag	UNP B1PNC0
E	11	ALA	-	expression tag	UNP B1PNC0
E	25	ILE	PHE	engineered mutation	UNP B1PNC0
E	35	GLN	ARG	engineered mutation	UNP B1PNC0
E	42	GLU	ASP	engineered mutation	UNP B1PNC0
E	55	ASP	ALA	engineered mutation	UNP B1PNC0
E	66	HIS	GLN	engineered mutation	UNP B1PNC0
E	67	VAL	ILE	engineered mutation	UNP B1PNC0
E	69	IO8	GLY	chromophore	UNP B1PNC0
E	?	-	TYR	chromophore	UNP B1PNC0
E	?	-	GLY	chromophore	UNP B1PNC0
E	77	TYR	PHE	engineered mutation	UNP B1PNC0
E	89	VAL	LYS	engineered mutation	UNP B1PNC0
E	110	VAL	SER	engineered mutation	UNP B1PNC0
E	125	ALA	PHE	engineered mutation	UNP B1PNC0
E	128	LYS	ILE	engineered mutation	UNP B1PNC0
E	146	VAL	ALA	engineered mutation	UNP B1PNC0
E	148	SER	TRP	engineered mutation	UNP B1PNC0

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Chain	Residue	Modelled	Actual	Comment	Reference
E	150	TRP	VAL	engineered mutation	UNP B1PNC0
E	151	SER	THR	engineered mutation	UNP B1PNC0
E	153	LYS	MET	engineered mutation	UNP B1PNC0
E	154	THR	LEU	engineered mutation	UNP B1PNC0
E	166	LYS	ASP	engineered mutation	UNP B1PNC0
E	168	SER	THR	engineered mutation	UNP B1PNC0
E	173	ASN	SER	engineered mutation	UNP B1PNC0
E	175	GLU	LYS	engineered mutation	UNP B1PNC0
E	178	ARG	GLN	engineered mutation	UNP B1PNC0
E	181	ALA	VAL	engineered mutation	UNP B1PNC0
E	184	THR	ASN	engineered mutation	UNP B1PNC0
E	195	TYR	ILE	engineered mutation	UNP B1PNC0
E	201	VAL	MET	engineered mutation	UNP B1PNC0
E	202	TYR	PHE	engineered mutation	UNP B1PNC0
E	213	GLU	LYS	engineered mutation	UNP B1PNC0
E	230	GLY	-	expression tag	UNP B1PNC0
E	231	MET	-	expression tag	UNP B1PNC0
E	232	ASP	-	expression tag	UNP B1PNC0
E	233	GLU	-	expression tag	UNP B1PNC0
E	234	LEU	-	expression tag	UNP B1PNC0
E	235	TYR	-	expression tag	UNP B1PNC0
E	236	LYS	-	expression tag	UNP B1PNC0
F	1	MET	-	initiating methionine	UNP B1PNC0
F	2	VAL	-	expression tag	UNP B1PNC0
F	3	SER	-	expression tag	UNP B1PNC0
F	4	LYS	-	expression tag	UNP B1PNC0
F	5	GLY	-	expression tag	UNP B1PNC0
F	6	GLU	-	expression tag	UNP B1PNC0
F	7	GLU	-	expression tag	UNP B1PNC0
F	8	ASP	-	expression tag	UNP B1PNC0
F	9	ASN	-	expression tag	UNP B1PNC0
F	10	MET	-	expression tag	UNP B1PNC0
F	11	ALA	-	expression tag	UNP B1PNC0
F	25	ILE	PHE	engineered mutation	UNP B1PNC0
F	35	GLN	ARG	engineered mutation	UNP B1PNC0
F	42	GLU	ASP	engineered mutation	UNP B1PNC0
F	55	ASP	ALA	engineered mutation	UNP B1PNC0
F	66	HIS	GLN	engineered mutation	UNP B1PNC0
F	67	VAL	ILE	engineered mutation	UNP B1PNC0
F	69	IO8	GLY	chromophore	UNP B1PNC0
F	?	-	TYR	chromophore	UNP B1PNC0
F	?	-	GLY	chromophore	UNP B1PNC0

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Chain	Residue	Modelled	Actual	Comment	Reference
F	77	TYR	PHE	engineered mutation	UNP B1PNC0
F	89	VAL	LYS	engineered mutation	UNP B1PNC0
F	110	VAL	SER	engineered mutation	UNP B1PNC0
F	125	ALA	PHE	engineered mutation	UNP B1PNC0
F	128	LYS	ILE	engineered mutation	UNP B1PNC0
F	146	VAL	ALA	engineered mutation	UNP B1PNC0
F	148	SER	TRP	engineered mutation	UNP B1PNC0
F	150	TRP	VAL	engineered mutation	UNP B1PNC0
F	151	SER	THR	engineered mutation	UNP B1PNC0
F	153	LYS	MET	engineered mutation	UNP B1PNC0
F	154	THR	LEU	engineered mutation	UNP B1PNC0
F	166	LYS	ASP	engineered mutation	UNP B1PNC0
F	168	SER	THR	engineered mutation	UNP B1PNC0
F	173	ASN	SER	engineered mutation	UNP B1PNC0
F	175	GLU	LYS	engineered mutation	UNP B1PNC0
F	178	ARG	GLN	engineered mutation	UNP B1PNC0
F	181	ALA	VAL	engineered mutation	UNP B1PNC0
F	184	THR	ASN	engineered mutation	UNP B1PNC0
F	195	TYR	ILE	engineered mutation	UNP B1PNC0
F	201	VAL	MET	engineered mutation	UNP B1PNC0
F	202	TYR	PHE	engineered mutation	UNP B1PNC0
F	213	GLU	LYS	engineered mutation	UNP B1PNC0
F	230	GLY	-	expression tag	UNP B1PNC0
F	231	MET	-	expression tag	UNP B1PNC0
F	232	ASP	-	expression tag	UNP B1PNC0
F	233	GLU	-	expression tag	UNP B1PNC0
F	234	LEU	-	expression tag	UNP B1PNC0
F	235	TYR	-	expression tag	UNP B1PNC0
F	236	LYS	-	expression tag	UNP B1PNC0
G	1	MET	-	initiating methionine	UNP B1PNC0
G	2	VAL	-	expression tag	UNP B1PNC0
G	3	SER	-	expression tag	UNP B1PNC0
G	4	LYS	-	expression tag	UNP B1PNC0
G	5	GLY	-	expression tag	UNP B1PNC0
G	6	GLU	-	expression tag	UNP B1PNC0
G	7	GLU	-	expression tag	UNP B1PNC0
G	8	ASP	-	expression tag	UNP B1PNC0
G	9	ASN	-	expression tag	UNP B1PNC0
G	10	MET	-	expression tag	UNP B1PNC0
G	11	ALA	-	expression tag	UNP B1PNC0
G	25	ILE	PHE	engineered mutation	UNP B1PNC0
G	35	GLN	ARG	engineered mutation	UNP B1PNC0

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Chain	Residue	Modelled	Actual	Comment	Reference
G	42	GLU	ASP	engineered mutation	UNP B1PNC0
G	55	ASP	ALA	engineered mutation	UNP B1PNC0
G	66	HIS	GLN	engineered mutation	UNP B1PNC0
G	67	VAL	ILE	engineered mutation	UNP B1PNC0
G	69	IO8	GLY	chromophore	UNP B1PNC0
G	?	-	TYR	chromophore	UNP B1PNC0
G	?	-	GLY	chromophore	UNP B1PNC0
G	77	TYR	PHE	engineered mutation	UNP B1PNC0
G	89	VAL	LYS	engineered mutation	UNP B1PNC0
G	110	VAL	SER	engineered mutation	UNP B1PNC0
G	125	ALA	PHE	engineered mutation	UNP B1PNC0
G	128	LYS	ILE	engineered mutation	UNP B1PNC0
G	146	VAL	ALA	engineered mutation	UNP B1PNC0
G	148	SER	TRP	engineered mutation	UNP B1PNC0
G	150	TRP	VAL	engineered mutation	UNP B1PNC0
G	151	SER	THR	engineered mutation	UNP B1PNC0
G	153	LYS	MET	engineered mutation	UNP B1PNC0
G	154	THR	LEU	engineered mutation	UNP B1PNC0
G	166	LYS	ASP	engineered mutation	UNP B1PNC0
G	168	SER	THR	engineered mutation	UNP B1PNC0
G	173	ASN	SER	engineered mutation	UNP B1PNC0
G	175	GLU	LYS	engineered mutation	UNP B1PNC0
G	178	ARG	GLN	engineered mutation	UNP B1PNC0
G	181	ALA	VAL	engineered mutation	UNP B1PNC0
G	184	THR	ASN	engineered mutation	UNP B1PNC0
G	195	TYR	ILE	engineered mutation	UNP B1PNC0
G	201	VAL	MET	engineered mutation	UNP B1PNC0
G	202	TYR	PHE	engineered mutation	UNP B1PNC0
G	213	GLU	LYS	engineered mutation	UNP B1PNC0
G	230	GLY	-	expression tag	UNP B1PNC0
G	231	MET	-	expression tag	UNP B1PNC0
G	232	ASP	-	expression tag	UNP B1PNC0
G	233	GLU	-	expression tag	UNP B1PNC0
G	234	LEU	-	expression tag	UNP B1PNC0
G	235	TYR	-	expression tag	UNP B1PNC0
G	236	LYS	-	expression tag	UNP B1PNC0
H	1	MET	-	initiating methionine	UNP B1PNC0
H	2	VAL	-	expression tag	UNP B1PNC0
H	3	SER	-	expression tag	UNP B1PNC0
H	4	LYS	-	expression tag	UNP B1PNC0
H	5	GLY	-	expression tag	UNP B1PNC0
H	6	GLU	-	expression tag	UNP B1PNC0

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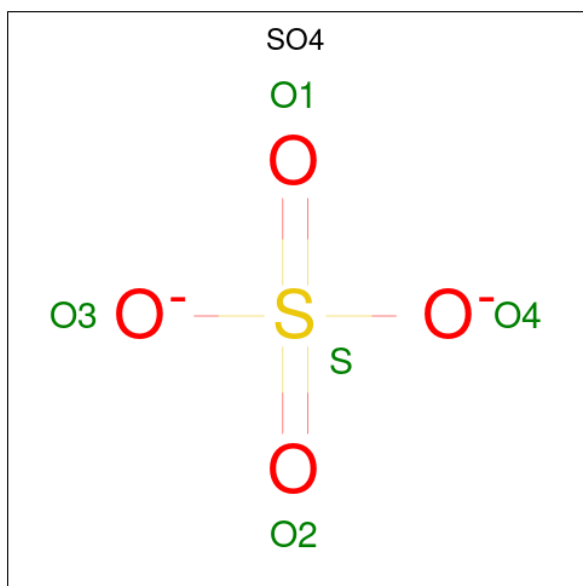
Chain	Residue	Modelled	Actual	Comment	Reference
H	7	GLU	-	expression tag	UNP B1PNC0
H	8	ASP	-	expression tag	UNP B1PNC0
H	9	ASN	-	expression tag	UNP B1PNC0
H	10	MET	-	expression tag	UNP B1PNC0
H	11	ALA	-	expression tag	UNP B1PNC0
H	25	ILE	PHE	engineered mutation	UNP B1PNC0
H	35	GLN	ARG	engineered mutation	UNP B1PNC0
H	42	GLU	ASP	engineered mutation	UNP B1PNC0
H	55	ASP	ALA	engineered mutation	UNP B1PNC0
H	66	HIS	GLN	engineered mutation	UNP B1PNC0
H	67	VAL	ILE	engineered mutation	UNP B1PNC0
H	69	IO8	GLY	chromophore	UNP B1PNC0
H	?	-	TYR	chromophore	UNP B1PNC0
H	?	-	GLY	chromophore	UNP B1PNC0
H	77	TYR	PHE	engineered mutation	UNP B1PNC0
H	89	VAL	LYS	engineered mutation	UNP B1PNC0
H	110	VAL	SER	engineered mutation	UNP B1PNC0
H	125	ALA	PHE	engineered mutation	UNP B1PNC0
H	128	LYS	ILE	engineered mutation	UNP B1PNC0
H	146	VAL	ALA	engineered mutation	UNP B1PNC0
H	148	SER	TRP	engineered mutation	UNP B1PNC0
H	150	TRP	VAL	engineered mutation	UNP B1PNC0
H	151	SER	THR	engineered mutation	UNP B1PNC0
H	153	LYS	MET	engineered mutation	UNP B1PNC0
H	154	THR	LEU	engineered mutation	UNP B1PNC0
H	166	LYS	ASP	engineered mutation	UNP B1PNC0
H	168	SER	THR	engineered mutation	UNP B1PNC0
H	173	ASN	SER	engineered mutation	UNP B1PNC0
H	175	GLU	LYS	engineered mutation	UNP B1PNC0
H	178	ARG	GLN	engineered mutation	UNP B1PNC0
H	181	ALA	VAL	engineered mutation	UNP B1PNC0
H	184	THR	ASN	engineered mutation	UNP B1PNC0
H	195	TYR	ILE	engineered mutation	UNP B1PNC0
H	201	VAL	MET	engineered mutation	UNP B1PNC0
H	202	TYR	PHE	engineered mutation	UNP B1PNC0
H	213	GLU	LYS	engineered mutation	UNP B1PNC0
H	230	GLY	-	expression tag	UNP B1PNC0
H	231	MET	-	expression tag	UNP B1PNC0
H	232	ASP	-	expression tag	UNP B1PNC0
H	233	GLU	-	expression tag	UNP B1PNC0
H	234	LEU	-	expression tag	UNP B1PNC0
H	235	TYR	-	expression tag	UNP B1PNC0

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Chain	Residue	Modelled	Actual	Comment	Reference
H	236	LYS	-	expression tag	UNP B1PNC0

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is water.

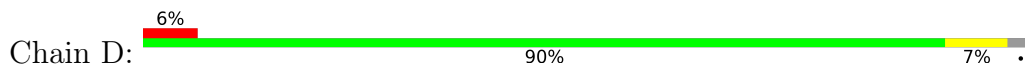
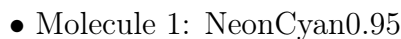
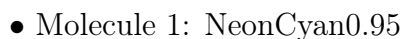
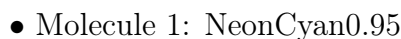
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	107	Total	O	0	0
			107	107		

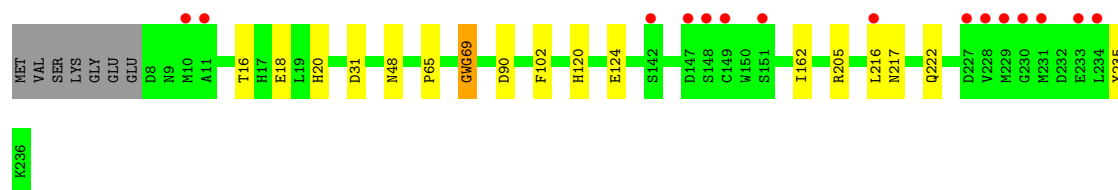
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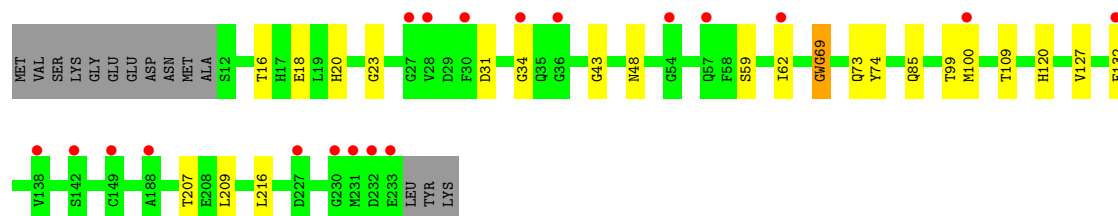
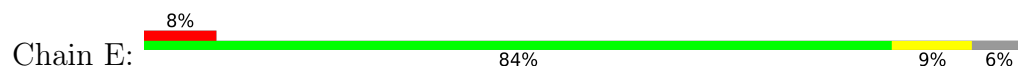
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	90	Total 90	O 90	0	0
3	C	78	Total 78	O 78	0	0
3	D	100	Total 100	O 100	0	1
3	E	33	Total 33	O 33	0	0
3	F	32	Total 32	O 32	0	0
3	G	23	Total 23	O 23	0	0
3	H	33	Total 33	O 33	0	0

- Molecule 1: NeonCyan0.95

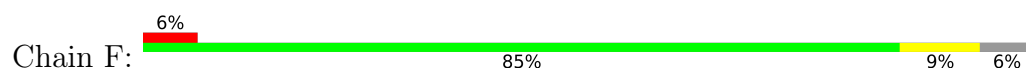




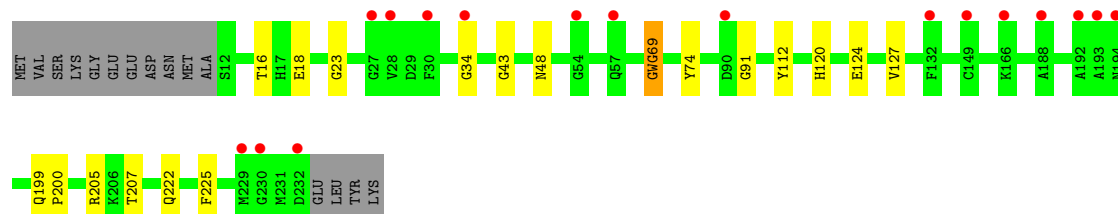
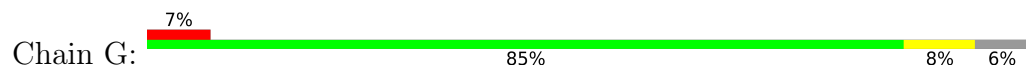
- Molecule 1: NeonCyan0.95



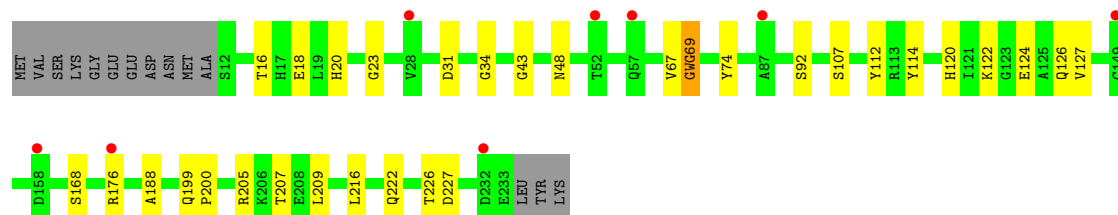
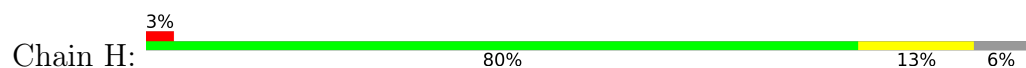
- Molecule 1: NeonCyan0.95



- Molecule 1: NeonCyan0.95



- Molecule 1: NeonCyan0.95



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	66.15Å 72.37Å 117.36Å 90.19° 90.04° 90.07°	Depositor
Resolution (Å)	48.90 – 1.95 48.90 – 1.95	Depositor EDS
% Data completeness (in resolution range)	96.0 (48.90-1.95) 76.8 (48.90-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.209 , 0.254 (Not available) , 0.272	Depositor DCC
R_{free} test set	7140 reflections (4.50%)	wwPDB-VP
Wilson B-factor (Å ²)	37.5	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 25.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.440 for h,-k,-l 0.438 for -h,k,-l 0.459 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14907	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.00% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: IO8, SO4

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	1.06	1/1884 (0.1%)	1.25	2/2553 (0.1%)
1	B	1.07	1/1854 (0.1%)	1.26	4/2514 (0.2%)
1	C	1.07	1/1891 (0.1%)	1.25	2/2563 (0.1%)
1	D	1.07	1/1883 (0.1%)	1.28	2/2554 (0.1%)
1	E	1.08	0/1781	1.27	0/2418
1	F	1.08	0/1811	1.25	0/2458
1	G	1.06	0/1790	1.23	0/2430
1	H	1.06	0/1790	1.26	1/2430 (0.0%)
All	All	1.07	4/14684 (0.0%)	1.26	11/19920 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	162	ILE	C-O	5.43	1.30	1.24
1	B	83	PRO	C-O	-5.26	1.17	1.24
1	C	73	GLN	C-O	5.20	1.30	1.24
1	D	162	ILE	C-O	5.04	1.29	1.24

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	67	VAL	CA-C-O	-5.83	110.89	120.80
1	B	102	PHE	CA-C-N	5.40	128.06	120.28
1	B	102	PHE	C-N-CA	5.40	128.06	120.28
1	C	102	PHE	CA-C-N	5.30	127.91	120.28
1	C	102	PHE	C-N-CA	5.30	127.91	120.28
1	B	165	PHE	CA-CB-CG	5.20	119.00	113.80
1	A	102	PHE	CA-C-N	5.17	127.72	120.28
1	A	102	PHE	C-N-CA	5.17	127.72	120.28
1	D	102	PHE	CA-C-N	5.10	127.62	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	102	PHE	C-N-CA	5.10	127.62	120.28
1	B	146	VAL	CA-C-O	-5.05	116.36	121.67

There are no chirality outliers.

There are no planarity outliers.

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	1842	0	1738	19	0
1	B	1813	0	1700	23	0
1	C	1849	0	1732	19	0
1	D	1841	0	1710	12	0
1	E	1749	0	1624	13	0
1	F	1770	0	1660	11	0
1	G	1752	0	1636	11	0
1	H	1755	0	1637	19	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
2	G	5	0	0	0	0
2	H	5	0	0	0	0
3	A	107	0	0	1	0
3	B	90	0	0	0	0
3	C	78	0	0	0	0
3	D	100	0	0	0	0
3	E	33	0	0	0	0
3	F	32	0	0	0	0
3	G	23	0	0	0	0
3	H	33	0	0	0	0
All	All	14907	0	13437	126	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 5.

All (126) 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:20:HIS:HE1	1:C:31:ASP:OD2	1.63	0.81
1:B:20:HIS:HE1	1:B:31:ASP:OD2	1.62	0.81
1:C:18:GLU:OE1	1:C:120:HIS:HE1	1.67	0.78
1:D:20:HIS:HE1	1:D:31:ASP:OD2	1.67	0.77
1:D:18:GLU:OE1	1:D:120:HIS:HE1	1.69	0.75
1:A:20:HIS:HE1	1:A:31:ASP:OD2	1.71	0.73
1:B:18:GLU:OE1	1:B:120:HIS:HE1	1.73	0.72
1:C:205:ARG:HG2	1:C:222:GLN:HB3	1.74	0.69
1:D:205:ARG:HE	1:D:222:GLN:HE21	1.41	0.68
1:A:18:GLU:OE1	1:A:120:HIS:HE1	1.77	0.66
1:D:16:THR:OG1	1:D:120:HIS:HD2	1.80	0.65
1:A:20:HIS:HD2	1:A:124:GLU:OE1	1.80	0.64
1:H:18:GLU:OE1	1:H:120:HIS:HE1	1.81	0.62
1:B:205:ARG:HG2	1:B:222:GLN:HB3	1.81	0.62
1:E:62:ILE:HD11	1:E:132:PHE:CZ	2.35	0.60
1:G:205:ARG:HG2	1:G:222:GLN:HG3	1.84	0.60
1:D:48[B]:ASN:OD1	1:D:217[B]:ASN:ND2	2.20	0.59
1:B:231:MET:HG3	1:B:231:MET:O	2.03	0.59
1:H:205:ARG:HG2	1:H:222:GLN:HB3	1.84	0.59
1:A:32:MET:HE1	1:A:216:LEU:HD13	1.87	0.57
1:A:16:THR:OG1	1:A:120:HIS:HD2	1.88	0.57
1:B:20:HIS:HD2	1:B:124:GLU:OE1	1.87	0.56
1:F:205:ARG:HH21	1:F:222:GLN:HE21	1.51	0.56
1:F:205:ARG:HG2	1:F:222:GLN:HG3	1.86	0.56
1:F:90:ASP:OD1	1:F:90:ASP:C	2.49	0.55
1:F:18:GLU:OE2	1:F:120:HIS:HE1	1.90	0.55
1:G:205:ARG:HH21	1:G:222:GLN:HE21	1.55	0.54
1:D:20:HIS:HD2	1:D:124:GLU:OE1	1.90	0.54
1:C:20:HIS:HD2	1:C:124:GLU:OE1	1.92	0.53
1:E:99:THR:HG22	1:E:109:THR:HG22	1.90	0.53
1:F:66:HIS:CE1	1:F:100[A]:MET:SD	3.02	0.53
1:B:34:GLY:HA3	1:B:48[B]:ASN:O	2.10	0.52
1:B:20:HIS:CE1	1:B:31:ASP:OD2	2.54	0.52
1:H:43:GLY:O	1:H:222:GLN:HG2	2.10	0.52
1:A:20:HIS:CD2	1:A:124:GLU:OE1	2.62	0.52
1:C:216:LEU:C	1:C:216:LEU:HD12	2.34	0.52
1:D:216:LEU:C	1:D:216:LEU:HD12	2.34	0.52
1:C:16:THR:OG1	1:C:120:HIS:HD2	1.93	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:THR:OG1	1:B:120:HIS:HD2	1.94	0.51
1:E:20:HIS:HE1	1:E:31:ASP:OD2	1.94	0.51
1:E:59:SER:O	1:E:62:ILE:HG22	2.10	0.51
1:A:228:VAL:HG12	1:A:231:MET:HG2	1.93	0.50
1:A:205:ARG:HG2	1:A:222:GLN:HB3	1.92	0.50
1:E:34:GLY:HA3	1:E:48:ASN:O	2.12	0.50
1:C:100[A]:MET:HG2	1:C:181:ALA:HB2	1.94	0.50
1:G:18:GLU:OE2	1:G:120:HIS:HE1	1.94	0.50
1:C:231:MET:O	1:C:231:MET:HG3	2.11	0.49
1:B:228:VAL:HG21	1:B:231:MET:HE3	1.95	0.49
1:B:78:PRO:HG2	1:B:226:THR:HG23	1.95	0.48
1:A:43:GLY:O	1:A:222:GLN:HG2	2.14	0.48
1:D:205:ARG:HE	1:D:222:GLN:NE2	2.08	0.47
1:G:43:GLY:HA2	1:G:74:TYR:O	2.13	0.47
1:A:231:MET:SD	1:A:234:LEU:HD23	2.55	0.47
1:E:18:GLU:OE1	1:E:120:HIS:HE1	1.98	0.47
1:B:34:GLY:HA3	1:B:48[A]:ASN:O	2.15	0.46
1:D:205:ARG:HG2	1:D:222:GLN:HG2	1.97	0.46
1:H:20:HIS:HE1	1:H:31:ASP:OD2	1.98	0.46
1:F:23:GLY:HA3	1:F:127:VAL:O	2.15	0.46
1:B:69:IO8:CZ3	1:B:207:THR:CG2	2.94	0.46
1:E:43:GLY:HA2	1:E:74:TYR:O	2.16	0.46
1:F:73:GLN:O	1:F:85:GLN:HG2	2.16	0.46
1:C:34:GLY:HA3	1:C:48[A]:ASN:O	2.16	0.46
1:A:175:GLU:HG2	3:A:505:HOH:O	2.15	0.46
1:H:209:LEU:HD11	1:H:216:LEU:HG	1.98	0.46
1:B:136:GLY:O	1:B:140:THR:HG23	2.16	0.46
1:H:16:THR:OG1	1:H:120:HIS:HD2	1.98	0.46
1:E:23:GLY:HA3	1:E:127:VAL:O	2.17	0.45
1:H:112:TYR:HA	1:H:124:GLU:O	2.15	0.45
1:A:78:PRO:HG2	1:A:226:THR:HG23	1.98	0.45
1:G:112:TYR:HA	1:G:124:GLU:O	2.16	0.45
1:H:20:HIS:CE1	1:H:31:ASP:OD2	2.70	0.45
1:B:100[A]:MET:HG2	1:B:181:ALA:HB2	1.98	0.45
1:D:20:HIS:CE1	1:D:31:ASP:OD2	2.58	0.45
1:H:226:THR:O	1:H:227:ASP:C	2.60	0.45
1:H:43:GLY:HA2	1:H:74:TYR:O	2.16	0.45
1:B:14:PRO:HB3	1:B:121:ILE:HD11	2.00	0.44
1:B:216:LEU:C	1:B:216:LEU:HD12	2.43	0.44
1:H:23:GLY:HA3	1:H:127:VAL:O	2.18	0.44
1:H:69:IO8:CH2	1:H:207:THR:CG2	2.96	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:82:SER:HB2	1:B:83:PRO:HD2	2.00	0.44
1:A:24:SER:HA	1:A:28:VAL:O	2.17	0.44
1:B:20:HIS:CD2	1:B:124:GLU:OE1	2.70	0.43
1:A:146:VAL:HG21	1:A:207:THR:CG2	2.48	0.43
1:E:69:IO8:CH2	1:E:207:THR:CG2	2.96	0.43
1:F:43:GLY:HA2	1:F:74:TYR:O	2.18	0.43
1:A:228:VAL:HG12	1:A:231:MET:CG	2.48	0.43
1:E:73:GLN:O	1:E:85:GLN:HG2	2.19	0.43
1:H:18:GLU:OE1	1:H:120:HIS:CE1	2.68	0.43
1:G:16:THR:OG1	1:G:120:HIS:HD2	2.01	0.43
1:C:24:SER:HA	1:C:28:VAL:O	2.19	0.43
1:C:49:LEU:HD12	1:C:216:LEU:HD11	2.01	0.43
1:G:34:GLY:HA3	1:G:48:ASN:O	2.19	0.43
1:C:20:HIS:CD2	1:C:124:GLU:OE1	2.71	0.42
1:B:209:LEU:HD11	1:B:216:LEU:HB2	2.00	0.42
1:B:231:MET:HE1	1:G:225:PHE:CZ	2.54	0.42
1:E:209:LEU:HD11	1:E:216:LEU:HG	2.01	0.42
1:A:215:GLU:C	1:A:216:LEU:HD12	2.45	0.42
1:B:69:IO8:CZ3	1:B:207:THR:HG21	2.48	0.42
1:C:20:HIS:CE1	1:C:31:ASP:OD2	2.55	0.42
1:C:216:LEU:HD12	1:C:216:LEU:O	2.19	0.42
1:D:18:GLU:OE1	1:D:120:HIS:CE1	2.60	0.42
1:H:92:SER:O	1:H:188:ALA:HB3	2.20	0.42
1:H:114:TYR:HA	1:H:122:LYS:O	2.20	0.42
1:C:62:ILE:HD12	1:C:62:ILE:HA	1.93	0.42
1:E:16:THR:OG1	1:E:120:HIS:HD2	2.02	0.42
1:H:34:GLY:HA3	1:H:48:ASN:O	2.19	0.42
1:C:34:GLY:HA3	1:C:48[B]:ASN:O	2.20	0.42
1:H:168:SER:OG	1:H:176:ARG:HD3	2.20	0.42
1:B:69:IO8:CH2	1:B:207:THR:HG23	2.50	0.41
1:G:23:GLY:HA3	1:G:127:VAL:O	2.20	0.41
1:E:20:HIS:CE1	1:E:31:ASP:OD2	2.73	0.41
1:F:96:VAL:HB	1:F:112:TYR:HB2	2.02	0.41
1:A:23:GLY:HA3	1:A:127:VAL:O	2.20	0.41
1:A:20:HIS:CE1	1:A:31:ASP:OD2	2.62	0.41
1:D:65:PRO:HB2	1:D:69:IO8:CD2	2.50	0.41
1:H:226:THR:HG22	1:H:227:ASP:N	2.34	0.41
1:F:147:ASP:OD2	1:F:168:SER:OG	2.39	0.41
1:C:18:GLU:OE1	1:C:120:HIS:CE1	2.59	0.41
1:F:16:THR:OG1	1:F:120:HIS:HD2	2.03	0.41
1:H:199:GLN:HA	1:H:200:PRO:C	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:209:LEU:HD11	1:C:216:LEU:HB2	2.04	0.40
1:G:69:IO8:CH2	1:G:207:THR:CG2	2.99	0.40
1:G:199:GLN:HA	1:G:200:PRO:C	2.46	0.40
1:A:147:ASP:OD2	1:A:168:SER:OG	2.35	0.40
1:C:65:PRO:HB2	1:C:69:IO8:CD2	2.51	0.40
1:B:152:LYS:O	1:B:163:SER:HA	2.22	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	227/234 (97%)	222 (98%)	5 (2%)	0	100	100
1	B	224/234 (96%)	217 (97%)	7 (3%)	0	100	100
1	C	228/234 (97%)	222 (97%)	6 (3%)	0	100	100
1	D	228/234 (97%)	221 (97%)	7 (3%)	0	100	100
1	E	216/234 (92%)	208 (96%)	8 (4%)	0	100	100
1	F	220/234 (94%)	210 (96%)	10 (4%)	0	100	100
1	G	217/234 (93%)	209 (96%)	7 (3%)	1 (0%)	24	16
1	H	217/234 (93%)	209 (96%)	8 (4%)	0	100	100
All	All	1777/1872 (95%)	1718 (97%)	58 (3%)	1 (0%)	48	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	91	GLY

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	197/200 (98%)	195 (99%)	2 (1%)	68	67
1	B	194/200 (97%)	193 (100%)	1 (0%)	81	82
1	C	197/200 (98%)	195 (99%)	2 (1%)	68	67
1	D	195/200 (98%)	193 (99%)	2 (1%)	68	67
1	E	184/200 (92%)	183 (100%)	1 (0%)	81	82
1	F	187/200 (94%)	183 (98%)	4 (2%)	47	41
1	G	185/200 (92%)	185 (100%)	0	100	100
1	H	184/200 (92%)	182 (99%)	2 (1%)	65	64
All	All	1523/1600 (95%)	1509 (99%)	14 (1%)	70	70

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	ASP
1	A	235	TYR
1	B	206	LYS
1	C	90	ASP
1	C	235	TYR
1	D	90	ASP
1	D	235	TYR
1	E	100	MET
1	F	57	GLN
1	F	62	ILE
1	F	90	ASP
1	F	234	LEU
1	H	107	SER
1	H	126	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	HIS
1	A	120	HIS
1	A	198	ASN
1	B	20	HIS
1	B	120	HIS
1	B	194	ASN
1	B	198	ASN
1	C	20	HIS
1	C	41	ASN
1	C	120	HIS
1	C	126	GLN
1	C	194	ASN
1	C	198	ASN
1	D	20	HIS
1	D	101	GLN
1	D	111	ASN
1	D	120	HIS
1	D	173	ASN
1	D	198	ASN
1	D	222	GLN
1	E	20	HIS
1	E	35	GLN
1	E	120	HIS
1	E	198	ASN
1	E	211	HIS
1	E	222	GLN
1	F	20	HIS
1	F	57	GLN
1	F	111	ASN
1	F	120	HIS
1	F	222	GLN
1	G	20	HIS
1	G	111	ASN
1	G	120	HIS
1	G	211	HIS
1	G	222	GLN
1	H	20	HIS
1	H	111	ASN
1	H	120	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	IO8	B	69	1	23,23,24	1.47	2 (8%)	29,32,34	3.00	13 (44%)
1	IO8	D	69	1	23,23,24	1.81	8 (34%)	29,32,34	3.54	11 (37%)
1	IO8	A	69	1	23,23,24	1.84	5 (21%)	29,32,34	2.92	12 (41%)
1	IO8	G	69	1	23,23,24	1.81	4 (17%)	29,32,34	2.94	9 (31%)
1	IO8	E	69	1	23,23,24	1.25	2 (8%)	29,32,34	3.47	15 (51%)
1	IO8	F	69	1	23,23,24	1.35	2 (8%)	29,32,34	3.22	11 (37%)
1	IO8	C	69	1	23,23,24	1.58	6 (26%)	29,32,34	3.06	14 (48%)
1	IO8	H	69	1	23,23,24	1.31	3 (13%)	29,32,34	2.68	10 (34%)

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	IO8	B	69	1	-	4/6/25/26	0/3/3/3
1	IO8	D	69	1	-	4/6/25/26	0/3/3/3
1	IO8	A	69	1	-	4/6/25/26	0/3/3/3
1	IO8	G	69	1	-	4/6/25/26	0/3/3/3
1	IO8	E	69	1	-	4/6/25/26	0/3/3/3
1	IO8	F	69	1	-	4/6/25/26	0/3/3/3
1	IO8	C	69	1	-	4/6/25/26	0/3/3/3
1	IO8	H	69	1	-	3/6/25/26	0/3/3/3

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	69	IO8	CB-CA2	5.63	1.49	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	69	IO8	CB-CG	-5.14	1.29	1.43
1	B	69	IO8	CB-CG	-3.95	1.33	1.43
1	D	69	IO8	CB-CG	-3.91	1.33	1.43
1	F	69	IO8	CB-CA2	3.65	1.45	1.38
1	A	69	IO8	CB-CA2	3.31	1.44	1.38
1	G	69	IO8	CA1-C1	-3.22	1.45	1.49
1	G	69	IO8	C2-N3	-3.18	1.32	1.40
1	D	69	IO8	CA2-C2	-3.15	1.45	1.48
1	B	69	IO8	C2-N3	-3.04	1.33	1.40
1	C	69	IO8	CE3-CD2	2.96	1.44	1.39
1	A	69	IO8	CE3-CD2	2.94	1.44	1.39
1	D	69	IO8	CE3-CD2	2.85	1.44	1.39
1	H	69	IO8	CB-CA2	2.83	1.44	1.38
1	C	69	IO8	C2-N3	-2.80	1.33	1.40
1	E	69	IO8	C1-N3	-2.79	1.32	1.37
1	C	69	IO8	CH2-CZ2	2.66	1.43	1.38
1	C	69	IO8	CZ3-CE3	2.65	1.43	1.38
1	E	69	IO8	CB-CA2	2.52	1.43	1.38
1	D	69	IO8	CH2-CZ2	2.47	1.43	1.38
1	A	69	IO8	C2-N3	-2.44	1.34	1.40
1	H	69	IO8	CA1-C1	-2.36	1.46	1.49
1	H	69	IO8	C2-N3	-2.35	1.34	1.40
1	G	69	IO8	C1-N3	-2.31	1.33	1.37
1	D	69	IO8	C2-N3	-2.28	1.34	1.40
1	A	69	IO8	CH2-CZ2	2.24	1.42	1.38
1	D	69	IO8	CZ2-CE2	2.22	1.43	1.39
1	C	69	IO8	CD2-CG	-2.06	1.42	1.45
1	D	69	IO8	CZ3-CE3	2.05	1.42	1.38
1	C	69	IO8	CA2-C2	-2.04	1.46	1.48
1	F	69	IO8	CB-CG	-2.02	1.38	1.43
1	D	69	IO8	CA1-N1	2.00	1.59	1.48

All (95) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	69	IO8	O2-C2-CA2	-11.45	123.72	131.02
1	F	69	IO8	O2-C2-CA2	-11.40	123.75	131.02
1	D	69	IO8	O2-C2-CA2	-9.39	125.03	131.02
1	G	69	IO8	O2-C2-CA2	-8.00	125.91	131.02
1	H	69	IO8	O2-C2-CA2	-7.51	126.22	131.02
1	A	69	IO8	O2-C2-CA2	-7.13	126.47	131.02
1	D	69	IO8	C1-CA1-N1	-7.05	97.26	112.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	69	IO8	O2-C2-CA2	-6.69	126.75	131.02
1	G	69	IO8	C1-CA1-N1	-6.65	98.14	112.85
1	B	69	IO8	C1-CA1-N1	-6.42	98.66	112.85
1	D	69	IO8	CG-CD1-NE1	-6.39	106.93	110.20
1	C	69	IO8	C2-N3-C1	-6.35	105.22	108.08
1	H	69	IO8	C1-CA1-N1	-6.31	98.90	112.85
1	C	69	IO8	C1-CA1-N1	-6.26	99.01	112.85
1	A	69	IO8	C1-CA1-N1	-6.20	99.13	112.85
1	D	69	IO8	CA2-C2-N3	6.17	108.68	103.50
1	B	69	IO8	CG-CD1-NE1	-6.08	107.10	110.20
1	F	69	IO8	CB-CA2-C2	6.05	130.59	122.47
1	F	69	IO8	C1-CA1-N1	-5.93	99.74	112.85
1	G	69	IO8	CB-CA2-C2	5.82	130.28	122.47
1	E	69	IO8	C2-CA2-N2	-5.73	104.84	108.95
1	G	69	IO8	C2-CA2-N2	-5.67	104.89	108.95
1	C	69	IO8	CA2-C2-N3	5.49	108.11	103.50
1	D	69	IO8	C2-CA2-N2	-5.32	105.14	108.95
1	B	69	IO8	O2-C2-CA2	-5.28	127.65	131.02
1	E	69	IO8	CB-CA2-C2	5.21	129.47	122.47
1	B	69	IO8	CA2-C2-N3	5.21	107.87	103.50
1	E	69	IO8	C1-CA1-N1	-5.08	101.61	112.85
1	D	69	IO8	C2-N3-C1	-5.07	105.80	108.08
1	D	69	IO8	CB-CA2-N2	5.03	135.15	124.60
1	A	69	IO8	CA2-C2-N3	4.98	107.69	103.50
1	E	69	IO8	CD2-CG-CB	-4.94	115.86	124.54
1	H	69	IO8	CB-CA2-C2	4.81	128.94	122.47
1	B	69	IO8	C2-N3-C1	-4.76	105.94	108.08
1	B	69	IO8	C2-CA2-N2	-4.76	105.54	108.95
1	D	69	IO8	CD2-CG-CD1	4.74	108.72	106.12
1	C	69	IO8	CB-CA2-N2	4.70	134.46	124.60
1	A	69	IO8	C2-N3-C1	-4.51	106.05	108.08
1	H	69	IO8	C2-CA2-N2	-4.38	105.81	108.95
1	A	69	IO8	C2-CA2-N2	-4.23	105.92	108.95
1	G	69	IO8	CA2-C2-N3	4.21	107.04	103.50
1	C	69	IO8	CG-CD1-NE1	-4.20	108.06	110.20
1	B	69	IO8	CD2-CG-CB	-4.08	117.37	124.54
1	A	69	IO8	CG-CD1-NE1	-4.08	108.12	110.20
1	E	69	IO8	CD1-CG-CB	4.07	137.88	129.59
1	H	69	IO8	CA2-C2-N3	3.93	106.80	103.50
1	F	69	IO8	CA3-N3-C2	3.86	132.16	123.67
1	A	69	IO8	C3-CA3-N3	3.83	121.14	112.43
1	E	69	IO8	CA3-N3-C2	3.71	131.83	123.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	69	IO8	CD2-CG-CD1	3.69	108.15	106.12
1	F	69	IO8	C2-CA2-N2	-3.63	106.35	108.95
1	C	69	IO8	N3-C1-N2	3.54	115.85	111.75
1	E	69	IO8	CA2-C2-N3	3.46	106.41	103.50
1	B	69	IO8	CB-CA2-N2	3.41	131.75	124.60
1	B	69	IO8	CD2-CG-CD1	3.34	107.96	106.12
1	A	69	IO8	CZ2-CE2-CD2	-3.30	119.00	122.19
1	E	69	IO8	CG-CD1-NE1	-3.25	108.54	110.20
1	C	69	IO8	C2-CA2-N2	-3.20	106.66	108.95
1	F	69	IO8	O2-C2-N3	3.13	130.84	124.31
1	C	69	IO8	CZ2-CE2-CD2	-3.04	119.25	122.19
1	E	69	IO8	CH2-CZ3-CE3	-3.03	116.50	120.24
1	B	69	IO8	C3-CA3-N3	3.01	119.27	112.43
1	F	69	IO8	CD2-CG-CB	-2.97	119.32	124.54
1	F	69	IO8	CG-CD1-NE1	-2.96	108.69	110.20
1	G	69	IO8	CG-CD1-NE1	-2.96	108.69	110.20
1	A	69	IO8	N3-C1-N2	2.95	115.16	111.75
1	C	69	IO8	C3-CA3-N3	2.94	119.11	112.43
1	G	69	IO8	CE2-CD2-CG	2.85	108.47	106.53
1	H	69	IO8	CG-CD1-NE1	-2.84	108.75	110.20
1	D	69	IO8	C3-CA3-N3	2.83	118.87	112.43
1	B	69	IO8	N3-C1-N2	2.77	114.95	111.75
1	E	69	IO8	CZ2-CE2-CD2	-2.69	119.58	122.19
1	C	69	IO8	CB-CA2-C2	-2.67	118.89	122.47
1	E	69	IO8	O2-C2-N3	2.67	129.88	124.31
1	A	69	IO8	CB-CA2-C2	2.64	126.02	122.47
1	D	69	IO8	CZ2-CE2-CD2	-2.63	119.64	122.19
1	H	69	IO8	C3-CA3-N3	2.55	118.24	112.43
1	G	69	IO8	CA3-N3-C2	2.54	129.24	123.67
1	B	69	IO8	CD1-CG-CB	2.49	134.67	129.59
1	A	69	IO8	CA3-N3-C2	2.46	129.08	123.67
1	C	69	IO8	CD2-CG-CD1	2.40	107.44	106.12
1	E	69	IO8	CE2-CD2-CG	2.39	108.16	106.53
1	F	69	IO8	CZ2-CE2-CD2	-2.35	119.91	122.19
1	F	69	IO8	CD1-CG-CB	2.35	134.37	129.59
1	E	69	IO8	CZ3-CE3-CD2	2.30	123.95	119.80
1	C	69	IO8	CA1-C1-N2	-2.29	121.20	124.28
1	F	69	IO8	CA2-C2-N3	2.28	105.42	103.50
1	B	69	IO8	O3-C3-CA3	-2.23	115.43	125.47
1	H	69	IO8	CA3-N3-C2	2.17	128.45	123.67
1	H	69	IO8	CE2-CD2-CG	2.17	108.01	106.53
1	E	69	IO8	C3-CA3-N3	2.10	117.20	112.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	69	IO8	CE2-NE1-CD1	2.09	110.93	109.08
1	D	69	IO8	CB-CA2-C2	-2.06	119.71	122.47
1	C	69	IO8	CH2-CZ2-CE2	2.02	122.80	118.69
1	H	69	IO8	O3-C3-CA3	-2.02	116.39	125.47

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	69	IO8	CA2-CB-CG-CD2
1	A	69	IO8	CA2-CB-CG-CD1
1	A	69	IO8	N2-CA2-CB-CG
1	B	69	IO8	CA2-CB-CG-CD2
1	B	69	IO8	CA2-CB-CG-CD1
1	B	69	IO8	N2-CA2-CB-CG
1	C	69	IO8	CA2-CB-CG-CD2
1	C	69	IO8	CA2-CB-CG-CD1
1	C	69	IO8	N2-CA2-CB-CG
1	D	69	IO8	CA2-CB-CG-CD2
1	D	69	IO8	CA2-CB-CG-CD1
1	D	69	IO8	N2-CA2-CB-CG
1	E	69	IO8	CA2-CB-CG-CD2
1	E	69	IO8	CA2-CB-CG-CD1
1	E	69	IO8	N2-CA2-CB-CG
1	F	69	IO8	CA2-CB-CG-CD2
1	F	69	IO8	CA2-CB-CG-CD1
1	F	69	IO8	N2-CA2-CB-CG
1	G	69	IO8	CA2-CB-CG-CD2
1	G	69	IO8	CA2-CB-CG-CD1
1	G	69	IO8	N2-CA2-CB-CG
1	H	69	IO8	CA2-CB-CG-CD2
1	H	69	IO8	CA2-CB-CG-CD1
1	H	69	IO8	N2-CA2-CB-CG
1	A	69	IO8	C2-CA2-CB-CG
1	B	69	IO8	C2-CA2-CB-CG
1	C	69	IO8	C2-CA2-CB-CG
1	D	69	IO8	C2-CA2-CB-CG
1	E	69	IO8	C2-CA2-CB-CG
1	F	69	IO8	C2-CA2-CB-CG
1	G	69	IO8	C2-CA2-CB-CG

There are no ring outliers.

6 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	69	IO8	3	0
1	D	69	IO8	1	0
1	G	69	IO8	1	0
1	E	69	IO8	1	0
1	C	69	IO8	1	0
1	H	69	IO8	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	SO4	A	301	-	4,4,4	0.48	0	6,6,6	0.32	0
2	SO4	B	301	-	4,4,4	0.56	0	6,6,6	0.20	0
2	SO4	C	301	-	4,4,4	0.45	0	6,6,6	0.36	0
2	SO4	D	301	-	4,4,4	0.42	0	6,6,6	0.33	0
2	SO4	H	301	-	4,4,4	0.42	0	6,6,6	0.19	0
2	SO4	G	301	-	4,4,4	0.48	0	6,6,6	0.25	0
2	SO4	F	301	-	4,4,4	0.43	0	6,6,6	0.16	0
2	SO4	E	301	-	4,4,4	0.36	0	6,6,6	0.19	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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	225/234 (96%)	0.36	10 (4%)	39	45	15, 34, 55, 81	6 (2%)
1	B	222/234 (94%)	0.45	12 (5%)	31	37	15, 34, 54, 88	6 (2%)
1	C	226/234 (96%)	0.42	9 (3%)	42	49	15, 35, 57, 89	6 (2%)
1	D	226/234 (96%)	0.49	15 (6%)	24	28	14, 35, 56, 92	6 (2%)
1	E	219/234 (93%)	0.77	19 (8%)	16	18	29, 46, 72, 90	1 (0%)
1	F	220/234 (94%)	0.73	13 (5%)	28	32	30, 45, 71, 83	4 (1%)
1	G	218/234 (93%)	0.82	17 (7%)	19	22	31, 47, 72, 84	3 (1%)
1	H	219/234 (93%)	0.73	8 (3%)	45	52	30, 47, 74, 85	2 (0%)
All	All	1775/1872 (94%)	0.59	103 (5%)	29	33	14, 41, 68, 92	34 (1%)

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	231	MET	5.1
1	B	232	ASP	4.7
1	D	234	LEU	4.1
1	D	148	SER	3.9
1	D	228	VAL	3.9
1	E	142	SER	3.9
1	F	158	ASP	3.8
1	D	230	GLY	3.8
1	H	158	ASP	3.5
1	C	228	VAL	3.4
1	F	233	GLU	3.4
1	E	231	MET	3.3
1	A	233	GLU	3.3
1	B	10	MET	3.3
1	F	234	LEU	3.2
1	A	10	MET	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	148	SER	3.1
1	E	149	CYS	3.1
1	E	232	ASP	3.1
1	G	132	PHE	2.9
1	E	62	ILE	2.9
1	C	234	LEU	2.9
1	G	229	MET	2.9
1	E	233	GLU	2.9
1	B	227	ASP	2.9
1	F	232	ASP	2.9
1	D	229	MET	2.9
1	C	148	SER	2.9
1	H	28	VAL	2.8
1	E	227	ASP	2.8
1	G	232	ASP	2.8
1	H	232	ASP	2.8
1	G	34	GLY	2.8
1	B	228	VAL	2.8
1	B	147	ASP	2.8
1	G	193	ALA	2.8
1	E	34	GLY	2.7
1	A	196	LEU	2.7
1	E	132	PHE	2.7
1	G	28	VAL	2.7
1	B	148	SER	2.7
1	C	10	MET	2.6
1	G	149	CYS	2.6
1	B	230	GLY	2.6
1	E	54	GLY	2.6
1	G	192	ALA	2.6
1	C	226	THR	2.5
1	E	57	GLN	2.5
1	F	57	GLN	2.5
1	C	8	ASP	2.5
1	E	30	PHE	2.5
1	F	52	THR	2.5
1	F	147	ASP	2.5
1	G	57	GLN	2.4
1	F	28	VAL	2.4
1	G	54	GLY	2.4
1	B	231	MET	2.4
1	B	206	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	G	90	ASP	2.4
1	E	230	GLY	2.4
1	H	57	GLN	2.4
1	D	147	ASP	2.3
1	D	227	ASP	2.3
1	F	196	LEU	2.3
1	D	11	ALA	2.3
1	A	149	CYS	2.3
1	D	216	LEU	2.3
1	G	166	LYS	2.3
1	D	233	GLU	2.3
1	A	228	VAL	2.3
1	G	194	ASN	2.3
1	A	234	LEU	2.2
1	C	196	LEU	2.2
1	A	232	ASP	2.2
1	C	147	ASP	2.2
1	B	229	MET	2.2
1	D	10	MET	2.2
1	D	149	CYS	2.2
1	D	142	SER	2.2
1	E	28	VAL	2.2
1	E	138	VAL	2.2
1	F	22	PHE	2.1
1	E	36	GLY	2.1
1	F	230	GLY	2.1
1	F	229	MET	2.1
1	E	188	ALA	2.1
1	G	188	ALA	2.1
1	G	30	PHE	2.1
1	A	235	TYR	2.1
1	A	147	ASP	2.1
1	B	142	SER	2.1
1	E	27	GLY	2.1
1	G	27	GLY	2.1
1	G	230	GLY	2.1
1	H	87	ALA	2.1
1	C	233	GLU	2.1
1	E	100	MET	2.0
1	H	52	THR	2.0
1	B	173	ASN	2.0
1	F	228	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
1	H	149	CYS	2.0
1	D	151	SER	2.0
1	H	176	ARG	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	IO8	A	69	21/22	0.93	0.12	24,30,40,43	0
1	IO8	C	69	21/22	0.94	0.12	23,28,38,41	0
1	IO8	H	69	21/22	0.94	0.10	34,40,44,49	0
1	IO8	D	69	21/22	0.95	0.08	23,29,41,42	0
1	IO8	E	69	21/22	0.95	0.08	34,40,45,48	0
1	IO8	F	69	21/22	0.95	0.10	33,40,44,48	0
1	IO8	B	69	21/22	0.95	0.08	25,31,38,44	0
1	IO8	G	69	21/22	0.96	0.08	35,41,46,47	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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(Å ²)	Q<0.9
2	SO4	H	301	5/5	0.94	0.10	49,52,54,59	0
2	SO4	F	301	5/5	0.95	0.09	46,49,51,59	0
2	SO4	G	301	5/5	0.97	0.07	48,51,56,59	0
2	SO4	C	301	5/5	0.97	0.09	35,37,44,56	0
2	SO4	E	301	5/5	0.98	0.05	45,47,53,54	0
2	SO4	B	301	5/5	0.98	0.07	34,39,44,52	0
2	SO4	A	301	5/5	0.98	0.08	36,36,45,50	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	D	301	5/5	0.98	0.07	37,39,45,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.