



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 11:58 AM UTC

PDB ID : 4ZFS / pdb_00004zfs
Title : Phototoxic Fluorescent Protein KillerOrange
Authors : Pletneva, N.V.; Pletnev, V.Z.; Pletnev, S.
Deposited on : 2015-04-21
Resolution : 2.01 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

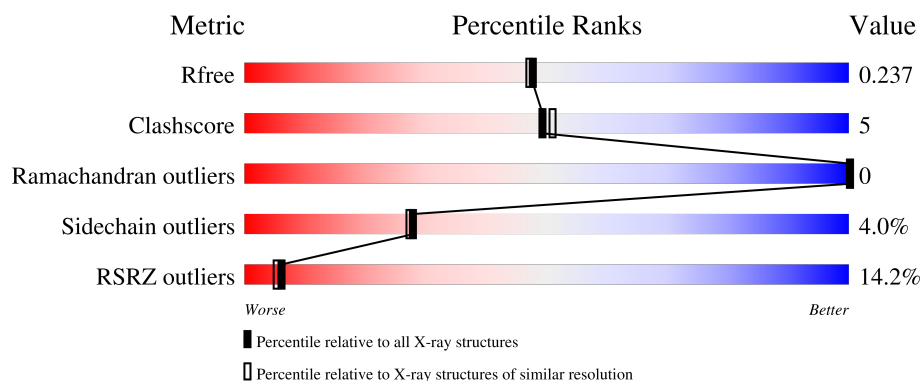
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	246	7% 79% 15% • 5%
1	B	246	14% 82% 10% •• 5%
1	C	246	11% 74% 16% • 7%
1	D	246	15% 78% 13% • 7%
1	E	246	19% 78% 13% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10068 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 KillerOrange.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	234	Total	C	N	O	S	0	4	0
			1870	1173	329	352	16			
1	B	234	Total	C	N	O	S	0	2	0
			1857	1166	328	347	16			
1	C	228	Total	C	N	O	S	0	6	0
			1843	1157	322	347	17			
1	D	228	Total	C	N	O	S	0	3	0
			1820	1144	319	341	16			
1	E	228	Total	C	N	O	S	0	2	0
			1816	1142	320	338	16			

There are 105 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	MET	-	initiating methionine	UNP Q2TCH5
A	-9	ARG	-	expression tag	UNP Q2TCH5
A	-8	GLY	-	expression tag	UNP Q2TCH5
A	-7	SER	-	expression tag	UNP Q2TCH5
A	-6	HIS	-	expression tag	UNP Q2TCH5
A	-5	HIS	-	expression tag	UNP Q2TCH5
A	-4	HIS	-	expression tag	UNP Q2TCH5
A	-3	HIS	-	expression tag	UNP Q2TCH5
A	-2	HIS	-	expression tag	UNP Q2TCH5
A	-1	HIS	-	expression tag	UNP Q2TCH5
A	0	GLY	-	expression tag	UNP Q2TCH5
A	1	SER	-	expression tag	UNP Q2TCH5
A	3	CYS	GLY	conflict	UNP Q2TCH5
A	65	4M9	GLN	chromophore	UNP Q2TCH5
A	65	4M9	TYR	chromophore	UNP Q2TCH5
A	65	4M9	GLY	chromophore	UNP Q2TCH5
A	113	SER	ASP	conflict	UNP Q2TCH5
A	145	SER	ASN	conflict	UNP Q2TCH5
A	177	LEU	PHE	conflict	UNP Q2TCH5

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Chain	Residue	Modelled	Actual	Comment	Reference
A	221	HIS	TYR	conflict	UNP Q2TCH5
A	236	GLN	GLU	conflict	UNP Q2TCH5
B	-10	MET	-	initiating methionine	UNP Q2TCH5
B	-9	ARG	-	expression tag	UNP Q2TCH5
B	-8	GLY	-	expression tag	UNP Q2TCH5
B	-7	SER	-	expression tag	UNP Q2TCH5
B	-6	HIS	-	expression tag	UNP Q2TCH5
B	-5	HIS	-	expression tag	UNP Q2TCH5
B	-4	HIS	-	expression tag	UNP Q2TCH5
B	-3	HIS	-	expression tag	UNP Q2TCH5
B	-2	HIS	-	expression tag	UNP Q2TCH5
B	-1	HIS	-	expression tag	UNP Q2TCH5
B	0	GLY	-	expression tag	UNP Q2TCH5
B	1	SER	-	expression tag	UNP Q2TCH5
B	3	CYS	GLY	conflict	UNP Q2TCH5
B	65	4M9	GLN	chromophore	UNP Q2TCH5
B	65	4M9	TYR	chromophore	UNP Q2TCH5
B	65	4M9	GLY	chromophore	UNP Q2TCH5
B	113	SER	ASP	conflict	UNP Q2TCH5
B	145	SER	ASN	conflict	UNP Q2TCH5
B	177	LEU	PHE	conflict	UNP Q2TCH5
B	221	HIS	TYR	conflict	UNP Q2TCH5
B	236	GLN	GLU	conflict	UNP Q2TCH5
C	-10	MET	-	initiating methionine	UNP Q2TCH5
C	-9	ARG	-	expression tag	UNP Q2TCH5
C	-8	GLY	-	expression tag	UNP Q2TCH5
C	-7	SER	-	expression tag	UNP Q2TCH5
C	-6	HIS	-	expression tag	UNP Q2TCH5
C	-5	HIS	-	expression tag	UNP Q2TCH5
C	-4	HIS	-	expression tag	UNP Q2TCH5
C	-3	HIS	-	expression tag	UNP Q2TCH5
C	-2	HIS	-	expression tag	UNP Q2TCH5
C	-1	HIS	-	expression tag	UNP Q2TCH5
C	0	GLY	-	expression tag	UNP Q2TCH5
C	1	SER	-	expression tag	UNP Q2TCH5
C	3	CYS	GLY	conflict	UNP Q2TCH5
C	65	4M9	GLN	chromophore	UNP Q2TCH5
C	65	4M9	TYR	chromophore	UNP Q2TCH5
C	65	4M9	GLY	chromophore	UNP Q2TCH5
C	113	SER	ASP	conflict	UNP Q2TCH5
C	145	SER	ASN	conflict	UNP Q2TCH5
C	177	LEU	PHE	conflict	UNP Q2TCH5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	221	HIS	TYR	conflict	UNP Q2TCH5
C	236	GLN	GLU	conflict	UNP Q2TCH5
D	-10	MET	-	initiating methionine	UNP Q2TCH5
D	-9	ARG	-	expression tag	UNP Q2TCH5
D	-8	GLY	-	expression tag	UNP Q2TCH5
D	-7	SER	-	expression tag	UNP Q2TCH5
D	-6	HIS	-	expression tag	UNP Q2TCH5
D	-5	HIS	-	expression tag	UNP Q2TCH5
D	-4	HIS	-	expression tag	UNP Q2TCH5
D	-3	HIS	-	expression tag	UNP Q2TCH5
D	-2	HIS	-	expression tag	UNP Q2TCH5
D	-1	HIS	-	expression tag	UNP Q2TCH5
D	0	GLY	-	expression tag	UNP Q2TCH5
D	1	SER	-	expression tag	UNP Q2TCH5
D	3	CYS	GLY	conflict	UNP Q2TCH5
D	65	4M9	GLN	chromophore	UNP Q2TCH5
D	65	4M9	TYR	chromophore	UNP Q2TCH5
D	65	4M9	GLY	chromophore	UNP Q2TCH5
D	113	SER	ASP	conflict	UNP Q2TCH5
D	145	SER	ASN	conflict	UNP Q2TCH5
D	177	LEU	PHE	conflict	UNP Q2TCH5
D	221	HIS	TYR	conflict	UNP Q2TCH5
D	236	GLN	GLU	conflict	UNP Q2TCH5
E	-10	MET	-	initiating methionine	UNP Q2TCH5
E	-9	ARG	-	expression tag	UNP Q2TCH5
E	-8	GLY	-	expression tag	UNP Q2TCH5
E	-7	SER	-	expression tag	UNP Q2TCH5
E	-6	HIS	-	expression tag	UNP Q2TCH5
E	-5	HIS	-	expression tag	UNP Q2TCH5
E	-4	HIS	-	expression tag	UNP Q2TCH5
E	-3	HIS	-	expression tag	UNP Q2TCH5
E	-2	HIS	-	expression tag	UNP Q2TCH5
E	-1	HIS	-	expression tag	UNP Q2TCH5
E	0	GLY	-	expression tag	UNP Q2TCH5
E	1	SER	-	expression tag	UNP Q2TCH5
E	3	CYS	GLY	conflict	UNP Q2TCH5
E	65	4M9	GLN	chromophore	UNP Q2TCH5
E	65	4M9	TYR	chromophore	UNP Q2TCH5
E	65	4M9	GLY	chromophore	UNP Q2TCH5
E	113	SER	ASP	conflict	UNP Q2TCH5
E	145	SER	ASN	conflict	UNP Q2TCH5
E	177	LEU	PHE	conflict	UNP Q2TCH5

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Chain	Residue	Modelled	Actual	Comment	Reference
E	221	HIS	TYR	conflict	UNP Q2TCH5
E	236	GLN	GLU	conflict	UNP Q2TCH5

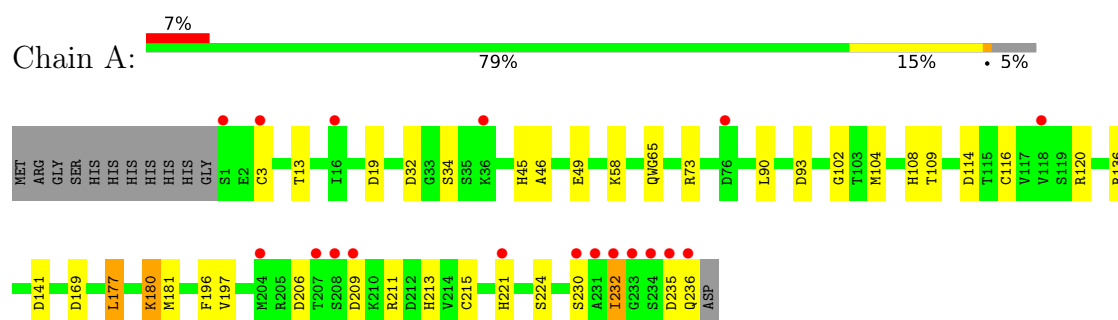
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	182	Total	O	0	0
			182	182		
2	B	185	Total	O	0	0
			185	185		
2	C	172	Total	O	0	0
			172	172		
2	D	167	Total	O	0	0
			167	167		
2	E	156	Total	O	0	0
			156	156		

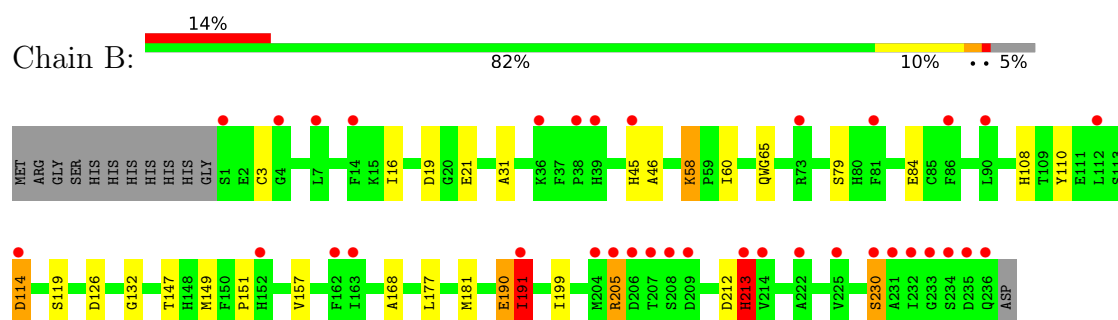
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

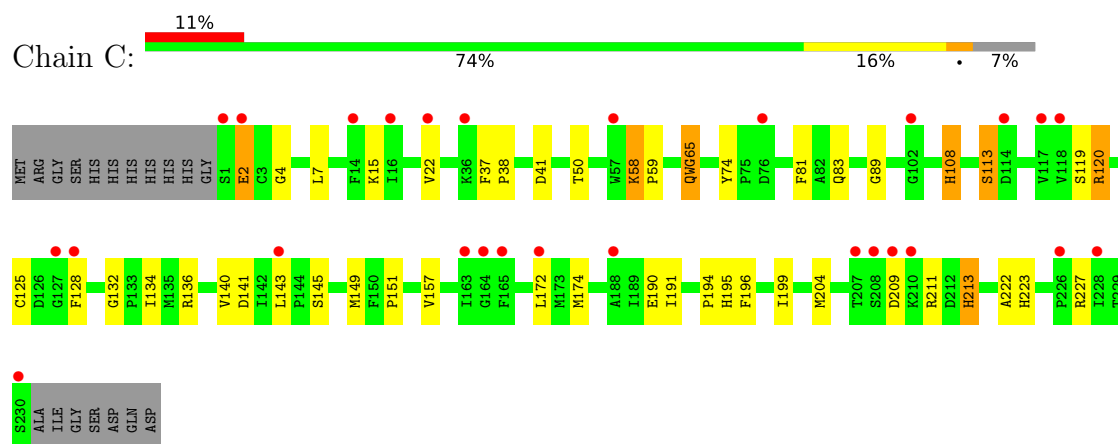
• Molecule 1: KillerOrange



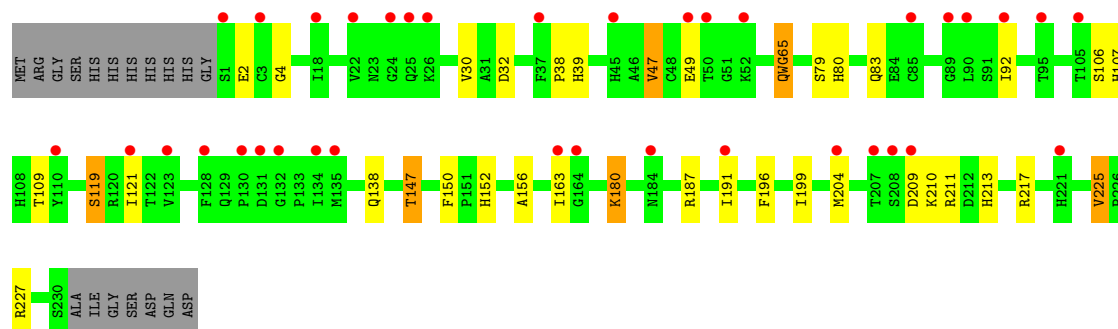
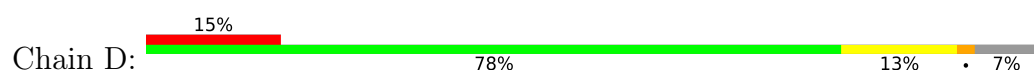
• Molecule 1: KillerOrange



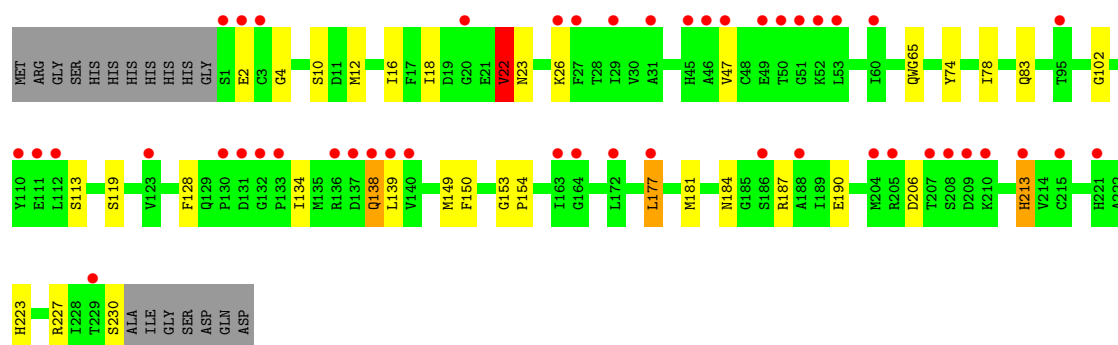
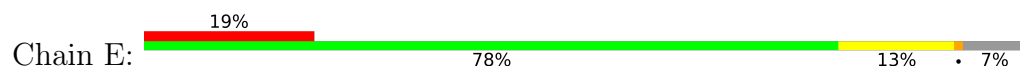
• Molecule 1: KillerOrange



• Molecule 1: KillerOrange



● Molecule 1: KillerOrange



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	128.93Å 202.06Å 116.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.85 – 2.01 29.85 – 2.01	Depositor EDS
% Data completeness (in resolution range)	99.4 (29.85-2.01) 99.6 (29.85-2.01)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.66 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.192 , 0.236 0.201 , 0.237	Depositor DCC
R_{free} test set	994 reflections (0.71%)	wwPDB-VP
Wilson B-factor (Å ²)	26.7	Xtriage
Anisotropy	1.253	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10068	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.88% 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: 4M9

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	1.35	8/1891 (0.4%)	1.15	1/2558 (0.0%)
1	B	1.37	7/1879 (0.4%)	1.22	12/2543 (0.5%)
1	C	1.29	8/1864 (0.4%)	1.16	5/2522 (0.2%)
1	D	1.28	4/1842 (0.2%)	1.16	3/2494 (0.1%)
1	E	1.24	4/1837 (0.2%)	1.16	7/2486 (0.3%)
All	All	1.31	31/9313 (0.3%)	1.17	28/12603 (0.2%)

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	191	ILE	N-CA	-8.32	1.39	1.46
1	B	191	ILE	CA-C	7.66	1.59	1.53
1	E	74	TYR	C-O	-7.10	1.15	1.24
1	B	168	ALA	C-O	-6.79	1.15	1.24
1	A	196	PHE	CA-C	6.53	1.60	1.52
1	C	195	HIS	C-O	6.45	1.30	1.24
1	B	60	ILE	CA-CB	6.10	1.61	1.54
1	C	223	HIS	C-O	6.09	1.31	1.23
1	C	41	ASP	C-O	-6.07	1.16	1.23
1	A	109	THR	CA-C	5.75	1.59	1.52
1	A	93	ASP	C-O	-5.71	1.17	1.24
1	D	80	HIS	CA-C	-5.64	1.46	1.53
1	B	58	LYS	C-O	-5.64	1.19	1.24
1	B	19	ASP	N-CA	-5.63	1.39	1.46
1	D	147	THR	C-O	5.61	1.31	1.23
1	C	81	PHE	C-O	5.61	1.30	1.24
1	E	78	ILE	C-O	-5.59	1.18	1.24
1	B	177	LEU	C-O	-5.56	1.17	1.23
1	A	197	VAL	C-O	5.49	1.30	1.24
1	D	106	SER	N-CA	-5.40	1.39	1.45
1	E	223	HIS	N-CA	-5.36	1.39	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	74	TYR	C-O	-5.35	1.17	1.24
1	A	177	LEU	CA-C	-5.34	1.46	1.52
1	C	174	MET	C-O	5.32	1.30	1.24
1	E	150	PHE	C-O	5.28	1.29	1.24
1	A	224	SER	C-O	-5.20	1.17	1.23
1	C	2	GLU	CA-C	5.18	1.59	1.53
1	C	151	PRO	C-O	-5.18	1.17	1.23
1	A	197	VAL	N-CA	5.08	1.53	1.46
1	A	102	GLY	C-O	-5.08	1.19	1.23
1	D	191	ILE	CA-CB	-5.07	1.48	1.53

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	132	GLY	CA-C-N	7.22	126.85	119.56
1	B	132	GLY	C-N-CA	7.22	126.85	119.56
1	E	74	TYR	CA-C-N	-6.75	113.76	120.98
1	E	74	TYR	C-N-CA	-6.75	113.76	120.98
1	B	213	HIS	N-CA-C	6.45	118.98	109.24
1	B	190	GLU	CA-C-N	-6.26	116.51	123.02
1	B	190	GLU	C-N-CA	-6.26	116.51	123.02
1	B	191	ILE	CB-CG1-CD1	-5.94	101.32	113.80
1	B	191	ILE	CA-C-N	5.78	125.69	119.85
1	B	191	ILE	C-N-CA	5.78	125.69	119.85
1	C	89	GLY	N-CA-C	5.72	119.25	112.33
1	C	213	HIS	N-CA-C	5.69	117.76	108.55
1	B	191	ILE	CA-CB-CG2	5.65	120.11	110.50
1	B	191	ILE	CB-CA-C	5.58	117.20	110.78
1	B	230	SER	N-CA-C	5.46	119.79	113.18
1	C	191	ILE	CB-CA-C	-5.45	104.63	110.95
1	B	3	CYS	N-CA-C	5.37	115.63	108.38
1	E	18	ILE	N-CA-C	5.36	115.57	107.80
1	E	22	VAL	CB-CA-C	-5.35	102.86	110.83
1	D	163	ILE	CB-CA-C	5.19	117.74	110.99
1	E	184	ASN	N-CA-C	5.13	117.77	111.82
1	A	3	CYS	N-CA-C	5.08	116.53	108.96
1	C	227	ARG	N-CA-C	-5.05	103.04	110.52
1	E	153	GLY	N-CA-C	-5.04	102.05	112.34
1	C	81	PHE	N-CA-C	5.04	117.42	111.33
1	E	177	LEU	CA-CB-CG	5.04	133.93	116.30
1	D	150	PHE	CA-C-N	5.03	124.78	119.76
1	D	150	PHE	C-N-CA	5.03	124.78	119.76

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	1870	0	1769	23	0
1	B	1857	0	1756	17	1
1	C	1843	0	1737	28	1
1	D	1820	0	1715	19	0
1	E	1816	0	1721	17	0
2	A	182	0	0	2	0
2	B	185	0	0	6	0
2	C	172	0	0	4	0
2	D	167	0	0	0	0
2	E	156	0	0	1	0
All	All	10068	0	8698	97	1

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 (97) 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:A:136[B]:ARG:HH11	1:A:136[B]:ARG:CG	1.63	1.11
1:A:136[B]:ARG:HG2	1:A:136[B]:ARG:NH1	1.49	0.96
1:C:2:GLU:HG2	1:C:7:LEU:HD21	1.66	0.77
1:A:32:ASP:OD2	1:C:113:SER:OG	2.02	0.76
1:C:2:GLU:CG	1:C:7:LEU:HD21	2.18	0.72
1:B:213:HIS:CD2	2:B:323:HOH:O	2.48	0.66
1:A:136[B]:ARG:HH11	1:A:136[B]:ARG:HG2	0.68	0.66
1:C:2:GLU:N	2:C:301:HOH:O	2.25	0.65
1:E:227[A]:ARG:NH1	2:E:301:HOH:O	2.29	0.65
1:B:45[A]:HIS:NE2	2:B:301:HOH:O	2.30	0.65
1:C:140:VAL:HG12	1:C:141:ASP:OD1	1.98	0.64
1:D:209:ASP:HB3	1:D:211:ARG:H	1.64	0.62
1:A:177:LEU:C	1:A:177:LEU:HD23	2.24	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:ARG:HB3	1:A:221:HIS:HD2	1.66	0.61
1:B:151:PRO:CB	1:B:191:ILE:HD11	2.31	0.60
1:C:136:ARG:HD3	2:C:416:HOH:O	2.01	0.60
1:D:196:PHE:HE2	1:D:225:VAL:HG13	1.66	0.60
1:E:149:MET:HE1	1:E:181:MET:HE1	1.85	0.58
1:C:149:MET:HE3	1:C:157:VAL:HG11	1.86	0.58
2:A:319:HOH:O	1:C:120:ARG:HG3	2.05	0.55
1:C:143:LEU:HD11	1:C:172:LEU:HD21	1.89	0.55
1:B:114:ASP:HA	2:B:428:HOH:O	2.05	0.55
1:C:50:THR:HG21	2:C:406:HOH:O	2.07	0.54
1:C:65:4M9:H10	1:C:65:4M9:O2	2.09	0.53
1:B:205:ARG:NH2	2:B:304:HOH:O	2.41	0.53
1:A:45:HIS:CD2	1:A:215:CYS:SG	3.03	0.52
1:B:151:PRO:HB3	1:B:191:ILE:HD11	1.92	0.51
1:D:209:ASP:HB2	1:D:213:HIS:NE2	2.25	0.51
1:C:2:GLU:HG3	1:C:7:LEU:HD21	1.93	0.51
1:B:149:MET:HE1	1:B:181:MET:HE2	1.92	0.51
1:D:4:GLY:HA2	1:D:83:GLN:O	2.10	0.51
1:A:46:ALA:O	1:A:213:HIS:HB2	2.12	0.50
1:A:235:ASP:O	1:A:236:GLN:CB	2.58	0.50
1:E:206:ASP:O	1:E:213:HIS:CD2	2.65	0.50
1:C:4:GLY:HA2	1:C:83:GLN:O	2.10	0.50
1:A:136[B]:ARG:CG	1:A:136[B]:ARG:NH1	2.36	0.50
1:C:211:ARG:O	1:C:213:HIS:CD2	2.66	0.49
1:A:235:ASP:O	1:A:236:GLN:HB2	2.13	0.48
1:B:230:SER:HB3	2:B:375:HOH:O	2.13	0.48
1:E:4:GLY:HA2	1:E:83:GLN:O	2.13	0.48
1:D:196:PHE:CE2	1:D:225:VAL:HG13	2.48	0.48
1:B:149:MET:HE3	1:B:157:VAL:HG11	1.96	0.47
1:D:79:SER:OG	1:E:190:GLU:OE1	2.24	0.47
1:D:227:ARG:NH2	1:E:187:ARG:O	2.48	0.47
1:A:49:GLU:HA	1:A:49:GLU:OE1	2.13	0.47
1:A:141:ASP:OD2	1:E:154:PRO:HG3	2.15	0.47
1:E:149:MET:HE1	1:E:181:MET:CE	2.45	0.47
1:B:16:ILE:HB	1:B:31:ALA:HB3	1.97	0.47
1:C:120:ARG:HD3	1:C:120:ARG:C	2.40	0.47
1:A:45:HIS:CD2	1:A:215:CYS:HG	2.34	0.46
1:B:110:TYR:CD2	1:B:119:SER:HB3	2.51	0.46
1:D:65:4M9:H10	1:D:65:4M9:O2	2.16	0.46
1:A:120:ARG:HD2	1:A:120:ARG:C	2.40	0.46
1:D:211:ARG:O	1:D:213:HIS:HD2	1.99	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:196:PHE:CE2	1:D:225:VAL:CG1	2.99	0.45
1:C:120:ARG:C	1:C:120:ARG:CD	2.89	0.45
1:C:132:GLY:O	1:C:136:ARG:HG3	2.16	0.45
1:E:22:VAL:O	1:E:23:ASN:C	2.60	0.45
1:C:22:VAL:HA	1:C:125:CYS:HB2	1.98	0.45
1:E:206:ASP:HB3	1:E:213:HIS:CD2	2.52	0.44
1:D:152:HIS:O	1:D:156:ALA:HB3	2.18	0.44
1:A:19[B]:ASP:OD2	1:C:120:ARG:NH1	2.51	0.44
1:A:13:THR:HA	1:A:34:SER:HA	2.00	0.44
1:E:10:SER:O	1:E:12:MET:HG3	2.17	0.44
1:C:58:LYS:CB	1:C:59:PRO:HD3	2.48	0.44
1:C:196:PHE:O	1:C:222:ALA:HA	2.18	0.44
1:A:90:LEU:HD11	1:A:181:MET:HB3	1.99	0.44
1:C:108:HIS:CD2	1:C:108:HIS:N	2.86	0.44
1:E:47:VAL:HG22	1:E:213:HIS:HB3	2.00	0.43
1:E:16:ILE:HG12	1:E:119:SER:OG	2.18	0.43
1:C:128:PHE:HE2	1:C:134:ILE:HD13	1.83	0.43
1:D:147:THR:HG21	1:D:199:ILE:HD12	2.01	0.43
1:A:116:CYS:SG	1:C:15:LYS:HE3	2.59	0.43
1:A:206:ASP:O	1:A:213:HIS:CD2	2.71	0.43
1:A:232:ILE:HG13	1:D:217:ARG:NH1	2.33	0.42
1:A:169:ASP:OD1	1:A:169:ASP:C	2.63	0.42
1:B:46:ALA:O	1:B:213:HIS:HB2	2.19	0.42
1:D:30:VAL:HG12	1:D:49:GLU:OE2	2.19	0.42
1:B:205:ARG:NH1	1:B:212:ASP:OD1	2.52	0.42
1:A:180:LYS:HE2	2:A:307:HOH:O	2.20	0.42
1:B:45[A]:HIS:CE1	2:B:301:HOH:O	2.71	0.42
1:D:109:THR:O	1:D:119:SER:HA	2.20	0.42
1:C:37:PHE:CD1	1:C:38:PRO:HA	2.56	0.41
1:C:199:ILE:HD13	1:C:199:ILE:HG21	1.83	0.41
1:E:102:GLY:HA3	1:E:128:PHE:CD1	2.56	0.41
1:D:38:PRO:O	1:D:39:HIS:HB2	2.20	0.41
1:D:107:HIS:O	1:D:121:ILE:HA	2.21	0.41
1:E:134:ILE:HG12	1:E:139:LEU:HD11	2.01	0.41
1:D:47:VAL:HB	1:D:213:HIS:HB3	2.03	0.41
1:D:92:ILE:HA	1:D:180:LYS:O	2.21	0.41
1:E:138[A]:GLN:HE21	1:E:138[A]:GLN:HB2	1.62	0.41
1:B:21:GLU:OE1	1:B:126:ASP:OD1	2.39	0.41
1:C:50:THR:CG2	2:C:406:HOH:O	2.66	0.41
1:C:149:MET:O	1:C:194:PRO:HA	2.21	0.40
1:B:147:THR:HG21	1:B:199:ILE:HD12	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:26:LYS:HB3	1:E:26:LYS:HE3	1.84	0.40
1:B:84:GLU:OE1	1:B:190:GLU:O	2.39	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:79:SER:OG	1:C:190[A]:GLU:OE2[6_445]	1.94	0.26

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	233/246 (95%)	224 (96%)	9 (4%)	0	100	100
1	B	231/246 (94%)	222 (96%)	9 (4%)	0	100	100
1	C	229/246 (93%)	221 (96%)	8 (4%)	0	100	100
1	D	226/246 (92%)	217 (96%)	9 (4%)	0	100	100
1	E	225/246 (92%)	216 (96%)	9 (4%)	0	100	100
All	All	1144/1230 (93%)	1100 (96%)	44 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	206/212 (97%)	197 (96%)	9 (4%)	25	24
1	B	204/212 (96%)	198 (97%)	6 (3%)	37	40
1	C	204/212 (96%)	196 (96%)	8 (4%)	28	28
1	D	201/212 (95%)	190 (94%)	11 (6%)	19	17
1	E	200/212 (94%)	192 (96%)	8 (4%)	28	27
All	All	1015/1060 (96%)	973 (96%)	42 (4%)	28	26

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	LYS
1	A	104	MET
1	A	108	HIS
1	A	114	ASP
1	A	180	LYS
1	A	209	ASP
1	A	211	ARG
1	A	230	SER
1	A	232	ILE
1	B	58	LYS
1	B	108	HIS
1	B	114	ASP
1	B	191	ILE
1	B	205	ARG
1	B	213	HIS
1	C	58	LYS
1	C	108	HIS
1	C	113	SER
1	C	119	SER
1	C	120	ARG
1	C	145	SER
1	C	204	MET
1	C	209	ASP
1	D	2	GLU
1	D	32[A]	ASP
1	D	32[B]	ASP
1	D	47	VAL
1	D	119	SER
1	D	138	GLN
1	D	180	LYS
1	D	187	ARG

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Mol	Chain	Res	Type
1	D	204	MET
1	D	210	LYS
1	D	225	VAL
1	E	2	GLU
1	E	22	VAL
1	E	113	SER
1	E	138[A]	GLN
1	E	138[B]	GLN
1	E	177	LEU
1	E	213	HIS
1	E	230	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	83	GLN
1	A	138	GLN
1	A	203	GLN
1	A	221	HIS
1	B	25	GLN
1	B	43	ASN
1	B	213	HIS
1	C	9	GLN
1	C	25	GLN
1	C	45	HIS
1	C	124	ASN
1	C	221	HIS
1	D	124	ASN
1	D	221	HIS
1	E	9	GLN
1	E	45	HIS
1	E	203	GLN
1	E	221	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

5 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	4M9	C	65	1	28,28,29	1.52	6 (21%)	32,39,41	2.70	12 (37%)
1	4M9	D	65	1	28,28,29	1.34	5 (17%)	32,39,41	2.26	11 (34%)
1	4M9	A	65	1	28,28,29	1.54	8 (28%)	32,39,41	2.66	11 (34%)
1	4M9	B	65	1	28,28,29	1.64	6 (21%)	32,39,41	2.67	15 (46%)
1	4M9	E	65	1	28,28,29	1.53	6 (21%)	32,39,41	3.06	10 (31%)

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	4M9	C	65	1	-	2/10/32/33	0/3/3/3
1	4M9	D	65	1	-	1/10/32/33	0/3/3/3
1	4M9	A	65	1	-	0/10/32/33	0/3/3/3
1	4M9	B	65	1	-	3/10/32/33	0/3/3/3
1	4M9	E	65	1	-	0/10/32/33	0/3/3/3

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	65	4M9	CA1-N1	3.89	1.36	1.27
1	B	65	4M9	C1-N2	-3.86	1.25	1.33
1	A	65	4M9	CA1-N1	3.76	1.35	1.27
1	C	65	4M9	CA1-N1	3.53	1.35	1.27
1	E	65	4M9	CA1-N1	3.47	1.35	1.27
1	E	65	4M9	CD1-CG2	-3.40	1.32	1.38
1	B	65	4M9	CA1-N1	3.39	1.35	1.27
1	E	65	4M9	CB2-CA2	-3.38	1.32	1.38
1	B	65	4M9	CB2-CA2	-3.26	1.32	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	65	4M9	C1-N2	-3.09	1.26	1.33
1	E	65	4M9	C1-N2	-3.06	1.26	1.33
1	C	65	4M9	CA2-C2	2.97	1.51	1.48
1	A	65	4M9	CD1-CG2	-2.90	1.33	1.38
1	A	65	4M9	CB2-CA2	-2.81	1.33	1.38
1	C	65	4M9	CB2-CA2	-2.80	1.33	1.38
1	C	65	4M9	C1-N2	-2.76	1.27	1.33
1	D	65	4M9	CD1-CG2	-2.69	1.33	1.38
1	B	65	4M9	CD2-CG2	-2.66	1.41	1.45
1	D	65	4M9	C1-N2	-2.50	1.28	1.33
1	A	65	4M9	CA3-N3	-2.42	1.42	1.47
1	E	65	4M9	CD1-NE1	2.38	1.41	1.37
1	B	65	4M9	CD2-CE2	-2.35	1.38	1.41
1	C	65	4M9	CD2-CG2	-2.30	1.42	1.45
1	E	65	4M9	CD2-CG2	-2.26	1.42	1.45
1	D	65	4M9	CD2-CE2	-2.21	1.38	1.41
1	A	65	4M9	CD2-CG2	-2.21	1.42	1.45
1	D	65	4M9	CB2-CA2	-2.19	1.34	1.38
1	A	65	4M9	CD1-NE1	2.18	1.40	1.37
1	C	65	4M9	CD1-CG2	-2.10	1.34	1.38
1	B	65	4M9	CD1-NE1	2.10	1.40	1.37
1	A	65	4M9	CD2-CE2	-2.07	1.38	1.41

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	65	4M9	O2-C2-CA2	9.22	136.90	131.02
1	E	65	4M9	CB2-CA2-C2	8.51	133.90	122.47
1	C	65	4M9	O2-C2-CA2	6.50	135.17	131.02
1	A	65	4M9	CB2-CA2-C2	6.39	131.05	122.47
1	D	65	4M9	CB2-CA2-C2	6.01	130.54	122.47
1	B	65	4M9	O2-C2-CA2	5.89	134.78	131.02
1	B	65	4M9	CB2-CA2-C2	5.86	130.34	122.47
1	E	65	4M9	CD2-CG2-CD1	5.55	109.17	106.12
1	C	65	4M9	CB2-CA2-C2	5.49	129.84	122.47
1	A	65	4M9	CA2-C2-N3	-5.47	98.91	103.50
1	A	65	4M9	O2-C2-CA2	5.15	134.31	131.02
1	C	65	4M9	C2-CA2-N2	4.88	112.45	108.95
1	A	65	4M9	CD2-CG2-CD1	4.83	108.78	106.12
1	B	65	4M9	CG-CD-N4	4.77	131.78	116.49
1	E	65	4M9	C3-CA3-N3	4.68	123.09	112.43
1	C	65	4M9	CA2-C2-N3	-4.63	99.61	103.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	65	4M9	CZ2-CE2-CD2	-4.50	117.83	122.19
1	E	65	4M9	CB2-CA2-N2	-4.26	115.65	124.60
1	D	65	4M9	CG-CB1-CA1	4.21	126.25	113.51
1	D	65	4M9	CZ2-CE2-CD2	-4.08	118.24	122.19
1	A	65	4M9	C2-CA2-N2	4.04	111.85	108.95
1	B	65	4M9	C3-CA3-N3	3.96	121.44	112.43
1	E	65	4M9	CZ2-CE2-CD2	-3.95	118.36	122.19
1	C	65	4M9	CD2-CG2-CD1	3.90	108.27	106.12
1	A	65	4M9	C3-CA3-N3	3.73	120.91	112.43
1	D	65	4M9	CD2-CG2-CD1	3.71	108.16	106.12
1	A	65	4M9	CB2-CA2-N2	-3.63	116.98	124.60
1	B	65	4M9	CE3-CD2-CE2	3.60	122.58	118.84
1	A	65	4M9	CE3-CD2-CE2	3.50	122.47	118.84
1	B	65	4M9	O5-CD-CG	-3.49	110.49	121.04
1	C	65	4M9	CE3-CD2-CE2	3.48	122.45	118.84
1	A	65	4M9	CZ2-CE2-CD2	-3.47	118.83	122.19
1	C	65	4M9	CB2-CA2-N2	-3.45	117.36	124.60
1	B	65	4M9	CG2-CD1-NE1	-3.41	108.45	110.20
1	B	65	4M9	CZ2-CE2-CD2	-3.41	118.89	122.19
1	B	65	4M9	CD2-CG2-CD1	3.26	107.91	106.12
1	B	65	4M9	N3-C1-N2	3.14	116.35	112.62
1	E	65	4M9	CA2-C2-N3	-3.11	100.89	103.50
1	C	65	4M9	N3-C1-N2	3.04	116.23	112.62
1	D	65	4M9	O5-CD-N4	-2.97	114.61	122.53
1	D	65	4M9	C3-CA3-N3	2.97	119.18	112.43
1	C	65	4M9	CD2-CG2-CB2	-2.95	119.36	124.54
1	D	65	4M9	CA2-C2-N3	-2.93	101.04	103.50
1	D	65	4M9	CG-CD-N4	2.88	125.74	116.49
1	B	65	4M9	CD2-CG2-CB2	-2.85	119.54	124.54
1	E	65	4M9	CE3-CD2-CE2	2.82	121.76	118.84
1	A	65	4M9	CE2-NE1-CD1	-2.65	106.72	109.08
1	B	65	4M9	CA2-C2-N3	-2.64	101.28	103.50
1	D	65	4M9	CE3-CD2-CE2	2.57	121.50	118.84
1	D	65	4M9	CB2-CA2-N2	-2.55	119.23	124.60
1	B	65	4M9	CB1-CG-CD	2.50	126.70	114.16
1	E	65	4M9	CG2-CD1-NE1	-2.47	108.94	110.20
1	B	65	4M9	CE3-CD2-CG2	-2.42	130.31	133.99
1	A	65	4M9	CD2-CG2-CB2	-2.25	120.58	124.54
1	C	65	4M9	O5-CD-CG	-2.25	114.24	121.04
1	B	65	4M9	CB2-CA2-N2	-2.23	119.91	124.60
1	C	65	4M9	O3-C3-CA3	-2.20	115.55	125.47
1	E	65	4M9	CE2-NE1-CD1	-2.18	107.14	109.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	65	4M9	O2-C2-CA2	2.11	132.37	131.02

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	C	65	4M9	C2-CA2-CB2-CG2
1	D	65	4M9	C1-CA1-CB1-CG
1	B	65	4M9	C1-CA1-CB1-CG
1	B	65	4M9	O5-CD-CG-CB1
1	B	65	4M9	N4-CD-CG-CB1
1	C	65	4M9	CA2-CB2-CG2-CD2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	65	4M9	1	0
1	D	65	4M9	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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	233/246 (94%)	0.81	18 (7%) 19 18	19, 38, 67, 117	4 (1%)
1	B	233/246 (94%)	0.97	35 (15%) 5 5	20, 40, 77, 112	2 (0%)
1	C	227/246 (92%)	1.02	27 (11%) 9 8	20, 42, 63, 115	6 (2%)
1	D	227/246 (92%)	1.23	36 (15%) 5 4	22, 45, 70, 142	3 (1%)
1	E	227/246 (92%)	1.22	47 (20%) 2 2	23, 45, 74, 114	2 (0%)
All	All	1147/1230 (93%)	1.05	163 (14%) 6 5	19, 42, 71, 142	17 (1%)

All (163) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	232	ILE	5.0
1	A	233	GLY	4.8
1	B	1	SER	4.5
1	A	232	ILE	4.4
1	A	234	SER	4.2
1	B	191	ILE	4.2
1	D	1	SER	4.1
1	A	231	ALA	3.9
1	E	207	THR	3.9
1	E	1	SER	3.8
1	E	52	LYS	3.7
1	A	235	ASP	3.6
1	E	210	LYS	3.6
1	B	235	ASP	3.5
1	C	226	PRO	3.4
1	D	45[A]	HIS	3.4
1	E	50	THR	3.3
1	D	131	ASP	3.3
1	E	138[A]	GLN	3.3
1	B	4	GLY	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	204	MET	3.3
1	A	209	ASP	3.3
1	C	209	ASP	3.3
1	A	207	THR	3.2
1	E	29	ILE	3.2
1	A	236	GLN	3.2
1	D	49	GLU	3.1
1	B	38	PRO	3.1
1	D	134	ILE	3.1
1	B	14	PHE	3.1
1	B	236	GLN	3.1
1	E	49	GLU	3.0
1	E	123	VAL	3.0
1	B	163	ILE	3.0
1	C	2	GLU	2.9
1	E	204	MET	2.9
1	D	164	GLY	2.9
1	A	1	SER	2.8
1	D	208	SER	2.8
1	E	221	HIS	2.8
1	B	222	ALA	2.8
1	D	22	VAL	2.8
1	E	47	VAL	2.8
1	E	132	GLY	2.8
1	B	90	LEU	2.8
1	C	207	THR	2.8
1	E	3	CYS	2.8
1	B	233	GLY	2.7
1	D	18	ILE	2.7
1	D	128	PHE	2.7
1	E	172	LEU	2.7
1	C	208	SER	2.7
1	C	57	TRP	2.7
1	B	207	THR	2.7
1	D	163	ILE	2.7
1	C	127	GLY	2.7
1	C	118	VAL	2.6
1	C	102	GLY	2.6
1	E	27	PHE	2.6
1	B	209	ASP	2.6
1	E	209	ASP	2.6
1	D	207	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	112	LEU	2.6
1	C	210	LYS	2.6
1	B	162	PHE	2.6
1	C	1	SER	2.6
1	D	132	GLY	2.6
1	E	213	HIS	2.6
1	E	53	LEU	2.6
1	C	163	ILE	2.5
1	E	177	LEU	2.5
1	B	225	VAL	2.5
1	E	131	ASP	2.5
1	B	208	SER	2.5
1	B	7	LEU	2.5
1	A	36	LYS	2.5
1	E	163	ILE	2.5
1	D	3	CYS	2.5
1	C	230	SER	2.5
1	E	31	ALA	2.5
1	D	52	LYS	2.5
1	E	140	VAL	2.4
1	B	114	ASP	2.4
1	C	114	ASP	2.4
1	A	3	CYS	2.4
1	D	110	TYR	2.4
1	B	231	ALA	2.4
1	B	36	LYS	2.4
1	B	45[A]	HIS	2.4
1	E	215	CYS	2.4
1	D	204	MET	2.4
1	E	46	ALA	2.4
1	E	139	LEU	2.4
1	D	37	PHE	2.4
1	C	76	ASP	2.4
1	D	191	ILE	2.4
1	B	73	ARG	2.4
1	C	128	PHE	2.4
1	D	26	LYS	2.4
1	E	208	SER	2.4
1	B	39	HIS	2.4
1	B	213	HIS	2.4
1	D	130	PRO	2.4
1	D	221	HIS	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	51	GLY	2.3
1	B	86	PHE	2.3
1	E	45	HIS	2.3
1	A	221	HIS	2.3
1	C	188	ALA	2.3
1	C	143	LEU	2.3
1	D	24	GLY	2.3
1	E	26	LYS	2.3
1	C	165	PHE	2.3
1	B	214	VAL	2.3
1	E	95	THR	2.3
1	E	186	SER	2.2
1	D	123	VAL	2.2
1	E	133	PRO	2.2
1	E	229	THR	2.2
1	E	137	ASP	2.2
1	E	112	LEU	2.2
1	C	14	PHE	2.2
1	D	95	THR	2.2
1	A	118	VAL	2.2
1	A	204	MET	2.2
1	B	206	ASP	2.2
1	D	209	ASP	2.2
1	C	164	GLY	2.2
1	E	111	GLU	2.2
1	E	110	TYR	2.2
1	D	184	ASN	2.2
1	C	16	ILE	2.2
1	D	92	ILE	2.2
1	E	60	ILE	2.2
1	D	25	GLN	2.2
1	D	90	LEU	2.2
1	E	20	GLY	2.1
1	E	188	ALA	2.1
1	B	205	ARG	2.1
1	D	85	CYS	2.1
1	D	50	THR	2.1
1	E	130	PRO	2.1
1	D	121	ILE	2.1
1	C	22	VAL	2.1
1	E	164	GLY	2.1
1	A	230	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	234	SER	2.1
1	D	105	THR	2.1
1	B	81	PHE	2.1
1	C	117	VAL	2.1
1	A	76	ASP	2.1
1	E	205	ARG	2.1
1	C	228	ILE	2.1
1	A	208	SER	2.1
1	B	230	SER	2.1
1	C	172	LEU	2.0
1	E	136	ARG	2.0
1	B	152	HIS	2.0
1	C	36[A]	LYS	2.0
1	D	89	GLY	2.0
1	A	16	ILE	2.0
1	E	2	GLU	2.0
1	D	135	MET	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	4M9	D	65	26/27	0.85	0.14	40,46,52,57	0
1	4M9	B	65	26/27	0.86	0.14	36,39,44,45	0
1	4M9	A	65	26/27	0.87	0.12	31,38,43,44	0
1	4M9	C	65	26/27	0.90	0.11	40,48,63,65	0
1	4M9	E	65	26/27	0.90	0.11	39,47,53,58	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.