



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 22, 2026 – 01:10 pm BST

PDB ID : 9T2C / pdb_00009t2c
Title : Human Single Domain TCR-like Antibody A10 in complex with NY-ESO-1(C165V):HLA-A0201
Authors : Cmunt, D.; Lau, K.; Coukos, G.; Dunn, S.M.
Deposited on : 2025-10-22
Resolution : 2.89 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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

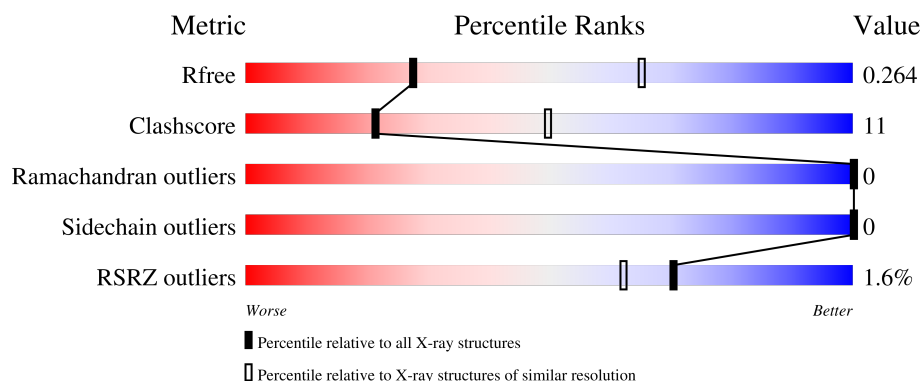
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	276	2% 73% 26% .
1	D	276	3% 72% 18% 11%
2	B	100	% 76% 23% .
2	E	100	% 79% 19% .
3	C	9	67% 33%

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Mol	Chain	Length	Quality of chain
3	F	9	67% 33%
4	G	147	% 58% 18% 24%
4	H	147	63% 18% •18%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 15041 atoms, of which 7292 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 MHC class I antigen.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	274	Total	C	H	N	O	S	0	0	0
			4232	1380	2033	396	414	9			
1	D	247	Total	C	H	N	O	S	0	0	0
			3870	1257	1869	361	375	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP A5I8L1
D	0	MET	-	initiating methionine	UNP A5I8L1

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	99	Total	C	H	N	O	S	0	0	0
			1577	518	768	134	153	4			
2	E	98	Total	C	H	N	O	S	0	0	0
			1595	521	779	138	154	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P61769
E	0	MET	-	initiating methionine	UNP P61769

- Molecule 3 is a protein called Cancer/testis antigen 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	9	Total	C	H	N	O	S	0	0	0
			158	51	83	11	12	1			
3	F	9	Total	C	H	N	O	S	0	0	0
			158	51	83	11	12	1			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	9	VAL	CYS	engineered mutation	UNP P78358
F	9	VAL	CYS	engineered mutation	UNP P78358

- Molecule 4 is a protein called A10 single-domain antibody.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
4	G	111	Total	C	H	N	O	S	0	0	0
			1657	537	807	142	169	2			
4	H	120	Total	C	H	N	O	S	0	0	0
			1789	578	870	153	186	2			

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Cl	0	0
			1	1		
5	D	1	Total	Cl	0	0
			1	1		

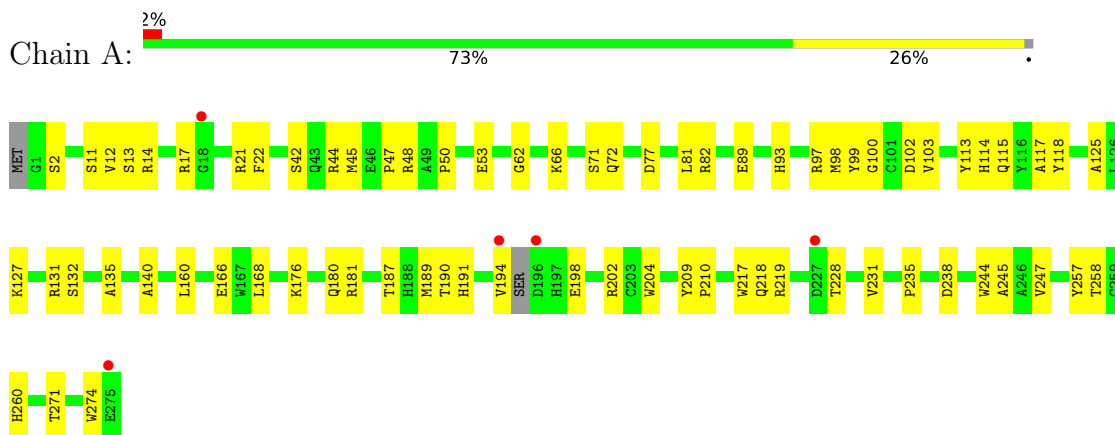
- Molecule 6 is water.

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

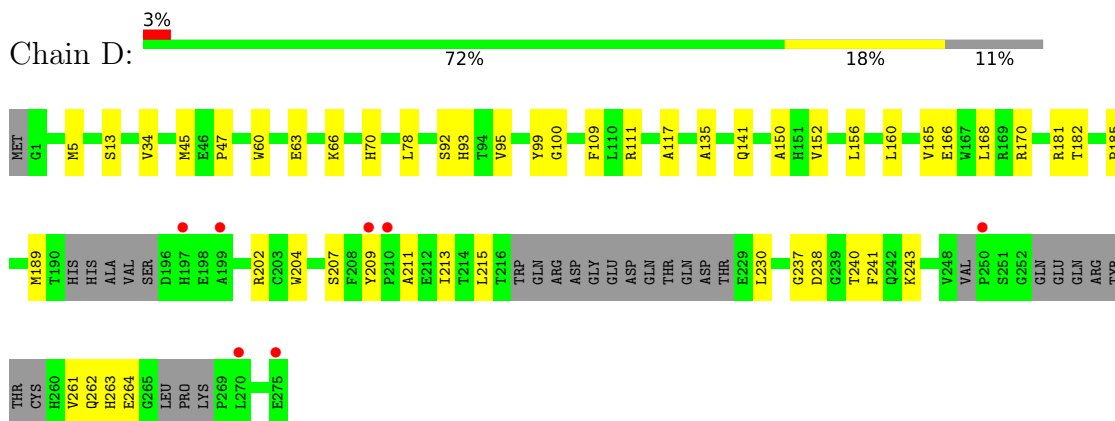
3 Residue-property plots [i](#)

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.

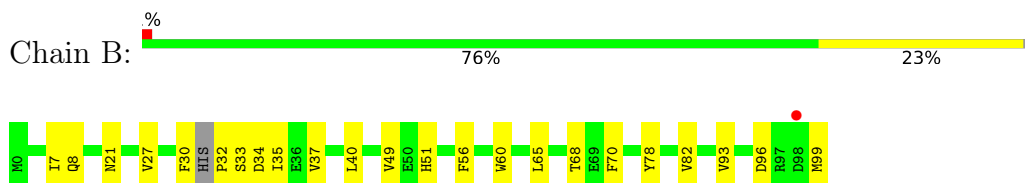
- Molecule 1: MHC class I antigen



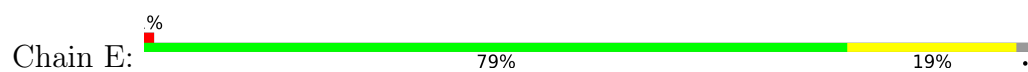
- Molecule 1: MHC class I antigen



- Molecule 2: Beta-2-microglobulin



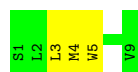
- Molecule 2: Beta-2-microglobulin



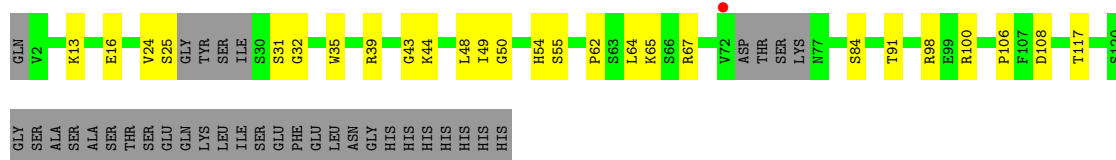
- Molecule 3: Cancer/testis antigen 1



- Molecule 3: Cancer/testis antigen 1



- Molecule 4: A10 single-domain antibody



- Molecule 4: A10 single-domain antibody



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.52Å 148.54Å 78.61Å 90.00° 104.98° 90.00°	Depositor
Resolution (Å)	19.68 – 2.89 19.68 – 2.89	Depositor EDS
% Data completeness (in resolution range)	99.1 (19.68-2.89) 98.8 (19.68-2.89)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.88Å)	Xtriage
Refinement program	PHENIX dev_5373	Depositor
R, R_{free}	0.180 , 0.264 0.180 , 0.264	Depositor DCC
R_{free} test set	1335 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	42.5	Xtriage
Anisotropy	0.287	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.44 , 59.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	15041	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.57% 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:
CL

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	0.59	0/2262	0.83	1/3074 (0.0%)
1	D	0.57	0/2055	0.80	0/2780
2	B	0.68	0/830	0.91	1/1123 (0.1%)
2	E	0.61	0/839	0.84	0/1137
3	C	0.95	0/76	1.12	0/103
3	F	0.62	0/76	0.88	0/103
4	G	0.62	0/871	0.85	0/1183
4	H	0.55	0/943	0.83	1/1284 (0.1%)
All	All	0.60	0/7952	0.84	3/10787 (0.0%)

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
2	E	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	56	PHE	CA-CB-CG	7.49	121.29	113.80
4	H	30	SER	N-CA-C	5.73	116.85	108.14
1	A	238	ASP	CA-CB-CG	5.15	117.75	112.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	E	97	ARG	Sidechain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2199	2033	2034	66	0
1	D	2001	1869	1869	37	0
2	B	809	768	768	19	0
2	E	816	779	779	13	0
3	C	75	83	83	4	0
3	F	75	83	83	9	0
4	G	850	807	805	18	0
4	H	919	870	870	22	0
5	A	1	0	0	0	0
5	D	1	0	0	0	0
6	A	2	0	0	1	0
6	D	1	0	0	0	0
All	All	7749	7292	7291	164	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 11.

All (164) 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:228:THR:HG22	1:A:247:VAL:CG1	1.75	1.15
1:A:14:ARG:NH2	1:A:21:ARG:HB2	1.61	1.13
1:A:14:ARG:HH22	1:A:21:ARG:HD3	1.42	0.84
4:G:13:LYS:O	4:G:16:GLU:HG3	1.82	0.79
1:D:5:MET:HB2	1:D:168:LEU:HD13	1.67	0.75
1:A:228:THR:HG22	1:A:247:VAL:HG12	1.66	0.75
1:A:228:THR:HG22	1:A:247:VAL:HG13	1.71	0.72
1:A:45:MET:CE	3:C:2:LEU:HD11	2.20	0.71
1:A:45:MET:HE1	3:C:2:LEU:HD11	1.72	0.71
1:A:14:ARG:HH21	1:A:21:ARG:HB2	1.56	0.70
2:B:27:VAL:HG11	2:B:35:ILE:HD13	1.72	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:ARG:HG2	1:A:17:ARG:CB	2.24	0.68
1:A:228:THR:HG22	1:A:247:VAL:HG11	1.73	0.67
4:H:35:TRP:HB2	4:H:52:ILE:HG23	1.79	0.65
1:A:77:ASP:OD1	3:C:9:VAL:HG22	1.99	0.63
1:A:82:ARG:HD2	1:A:89:GLU:HA	1.81	0.63
4:G:43:GLY:O	4:G:44:LYS:HD2	1.99	0.62
4:H:72:VAL:HG23	4:H:79:PHE:HB3	1.82	0.61
1:D:117:ALA:HB2	2:E:60:TRP:CE2	2.36	0.61
1:A:14:ARG:HH22	1:A:21:ARG:HB2	1.62	0.60
1:D:211:ALA:HA	1:D:241:PHE:CD2	2.36	0.60
4:H:100:ARG:HH11	4:H:100:ARG:HG3	1.64	0.60
2:E:1:ILE:HD12	2:E:1:ILE:O	2.01	0.60
1:D:109:PHE:HD1	1:D:165:VAL:HG11	1.67	0.59
1:A:14:ARG:HH22	1:A:21:ARG:CD	2.14	0.59
1:A:102:ASP:OD2	1:A:113:TYR:OH	2.10	0.59
1:D:207:SER:HA	1:D:240:THR:HB	1.82	0.59
2:E:41:LYS:NZ	2:E:78:TYR:OH	2.34	0.58
1:D:211:ALA:HB2	1:D:241:PHE:HE2	1.69	0.57
2:E:1:ILE:HD12	2:E:1:ILE:C	2.29	0.57
1:A:11:SER:HA	1:A:21:ARG:O	2.05	0.57
4:H:14:PRO:O	4:H:15:SER:OG	2.20	0.57
1:A:189:MET:SD	1:A:217:TRP:HH2	2.29	0.56
1:A:81:LEU:HD13	1:A:118:TYR:CD1	2.40	0.56
1:A:166:GLU:OE2	4:G:100:ARG:NH2	2.38	0.56
1:A:98:MET:HG2	1:A:99:TYR:N	2.21	0.55
1:D:150:ALA:HB3	1:D:152:VAL:HG23	1.89	0.55
2:B:7:ILE:HG21	2:B:93:VAL:HG21	1.88	0.54
1:A:218:GLN:OE1	1:A:260:HIS:NE2	2.41	0.53
4:H:13:LYS:HB2	4:H:16:GLU:CD	2.34	0.53
1:A:218:GLN:O	1:A:257:TYR:HA	2.08	0.53
4:G:24:VAL:HG12	4:G:25:SER:N	2.23	0.53
4:G:91:THR:HG23	4:G:117:THR:HA	1.90	0.53
1:A:209:TYR:HB3	1:A:210:PRO:HD3	1.92	0.52
1:A:228:THR:HA	1:A:247:VAL:HG12	1.91	0.52
2:B:21:ASN:HB3	2:B:70:PHE:CE1	2.44	0.52
4:G:54:HIS:O	4:G:55:SER:CB	2.57	0.51
1:A:228:THR:CG2	1:A:247:VAL:HG12	2.38	0.51
1:A:14:ARG:NH2	1:A:21:ARG:HD3	2.19	0.51
2:B:7:ILE:CG2	2:B:93:VAL:HG21	2.41	0.51
1:D:230:LEU:HD21	1:D:243:LYS:CE	2.40	0.51
1:A:176:LYS:HA	1:A:180:GLN:HG3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:35:TRP:CH2	4:G:98:ARG:HG3	2.46	0.50
1:D:189:MET:HA	1:D:202:ARG:O	2.12	0.50
1:D:135:ALA:HB3	1:D:141:GLN:HG2	1.92	0.50
1:D:238:ASP:HB3	2:E:12:ARG:NH1	2.25	0.50
1:A:258:THR:HG21	1:A:260:HIS:CE1	2.47	0.49
1:D:63:GLU:OE2	4:H:103:SER:OG	2.30	0.49
1:D:78:LEU:CD2	1:D:95:VAL:HG23	2.43	0.49
4:G:31:SER:O	4:G:54:HIS:ND1	2.45	0.49
1:A:115:GLN:HG2	1:A:125:ALA:CB	2.43	0.49
1:A:115:GLN:HG2	1:A:125:ALA:HB1	1.95	0.49
4:H:30:SER:OG	4:H:31:SER:N	2.46	0.49
1:D:13:SER:HB2	1:D:93:HIS:H	1.78	0.48
1:A:235:PRO:HG2	2:B:65:LEU:HD13	1.96	0.48
2:B:30:PHE:O	2:B:32:PRO:N	2.46	0.48
1:D:34:VAL:HB	1:D:45:MET:HE2	1.94	0.48
1:A:219:ARG:CB	1:A:257:TYR:CE1	2.96	0.48
1:A:13:SER:HB2	1:A:93:HIS:H	1.78	0.48
3:F:5:TRP:HE1	4:H:51:SER:HB3	1.79	0.48
4:G:67:ARG:HD2	4:G:84:SER:O	2.14	0.48
1:D:185:PRO:HD3	1:D:263:HIS:CD2	2.49	0.48
1:A:194:VAL:HB	1:A:198:GLU:HB2	1.96	0.47
1:A:231:VAL:HG21	2:B:8:GLN:OE1	2.14	0.47
1:D:100:GLY:O	1:D:160:LEU:HD22	2.14	0.47
1:A:62:GLY:O	1:A:66:LYS:HG3	2.15	0.47
1:A:66:LYS:HD3	3:C:4:MET:HA	1.95	0.47
4:H:58:THR:O	4:H:58:THR:HG23	2.13	0.47
1:A:72:GLN:HG2	4:G:62:PRO:HG3	1.97	0.47
1:A:135:ALA:HB1	1:A:140:ALA:HB3	1.96	0.47
1:D:135:ALA:CB	1:D:141:GLN:HG2	2.44	0.47
1:D:66:LYS:CD	3:F:4:MET:HA	2.44	0.47
4:G:39:ARG:HG2	4:G:49:ILE:HD11	1.95	0.47
1:A:127:LYS:HD3	1:A:132:SER:HB2	1.97	0.47
1:D:230:LEU:HD21	1:D:243:LYS:HE2	1.97	0.47
1:A:22:PHE:CD2	1:A:71:SER:HB2	2.49	0.47
1:A:135:ALA:HB1	1:A:140:ALA:CB	2.45	0.47
4:G:48:LEU:HD11	4:G:50:GLY:O	2.14	0.47
1:A:191:HIS:CB	1:A:274:TRP:CZ2	2.98	0.46
2:B:51:HIS:HA	2:B:65:LEU:O	2.15	0.46
3:F:4:MET:HE2	4:H:34:TYR:CD1	2.51	0.46
4:G:24:VAL:CG1	4:G:25:SER:N	2.78	0.46
4:G:39:ARG:CG	4:G:49:ILE:HD11	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:HIS:HB3	1:A:274:TRP:CZ2	2.51	0.46
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.51	0.46
1:D:261:VAL:C	1:D:262:GLN:HG2	2.41	0.46
4:G:64:LEU:O	4:G:65:LYS:C	2.58	0.45
1:A:100:GLY:O	1:A:160:LEU:HD22	2.16	0.45
4:H:72:VAL:O	4:H:72:VAL:HG13	2.15	0.45
2:B:27:VAL:HG11	2:B:35:ILE:CD1	2.42	0.45
1:A:131:ARG:NH2	6:A:401:HOH:O	2.50	0.45
4:H:39:ARG:HB3	4:H:49:ILE:HD11	1.98	0.45
1:A:50:PRO:HA	1:A:53:GLU:OE2	2.17	0.45
1:D:156:LEU:CD2	3:F:3:LEU:HD21	2.47	0.45
2:E:6:LYS:HE3	2:E:29:GLY:HA3	1.99	0.44
3:F:5:TRP:NE1	4:H:51:SER:HB3	2.33	0.44
1:D:156:LEU:HD23	3:F:3:LEU:HD21	1.99	0.44
2:E:41:LYS:O	2:E:42:ASN:C	2.61	0.44
1:A:187:THR:HA	1:A:204:TRP:O	2.17	0.44
1:D:237:GLY:O	2:E:12:ARG:NH2	2.51	0.44
2:E:31:HIS:CD2	2:E:62:PHE:CE2	3.05	0.44
2:B:96:ASP:OD1	2:B:96:ASP:C	2.59	0.44
2:B:96:ASP:O	2:B:99:MET:HG3	2.18	0.44
1:A:189:MET:HG3	1:A:190:THR:N	2.32	0.43
3:F:4:MET:CE	4:H:34:TYR:CD1	3.01	0.43
1:A:97:ARG:HH11	1:A:114:HIS:CE1	2.37	0.43
1:A:127:LYS:HB3	1:A:127:LYS:HE3	1.73	0.43
1:D:13:SER:HB2	1:D:92:SER:HA	2.00	0.43
2:E:27:VAL:HG23	2:E:30:PHE:HE2	1.82	0.43
1:D:70:HIS:NE2	1:D:99:TYR:OH	2.42	0.43
1:D:109:PHE:CZ	1:D:111:ARG:C	2.96	0.43
2:B:40:LEU:O	2:B:78:TYR:HA	2.18	0.43
1:D:213:ILE:HG13	1:D:263:HIS:HB2	1.99	0.43
4:H:48:LEU:HG	4:H:49:ILE:N	2.33	0.43
1:A:181:ARG:NH1	1:A:209:TYR:HB2	2.34	0.43
4:G:31:SER:OG	4:G:32:GLY:N	2.51	0.43
2:E:54:LEU:HA	2:E:64:LEU:HD21	2.00	0.43
2:B:33:SER:O	2:B:34:ASP:HB2	2.19	0.42
2:B:37:VAL:HG22	2:B:82:VAL:HG22	2.00	0.42
2:B:49:VAL:HG22	2:B:68:THR:HB	2.00	0.42
1:A:166:GLU:CD	4:G:100:ARG:HH21	2.27	0.42
1:A:2:SER:HB2	1:A:103:VAL:O	2.19	0.42
1:A:271:THR:HG23	1:A:271:THR:O	2.20	0.42
1:D:182:THR:O	1:D:182:THR:HG22	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:MET:HE3	1:A:99:TYR:HA	2.02	0.42
1:A:209:TYR:O	1:A:210:PRO:C	2.61	0.42
1:D:211:ALA:HA	1:D:241:PHE:HD2	1.82	0.42
4:H:53:TYR:CG	4:H:54:HIS:N	2.87	0.42
1:A:12:VAL:HG12	1:A:13:SER:O	2.19	0.41
1:A:47:PRO:O	1:A:48:ARG:HG2	2.19	0.41
1:D:263:HIS:O	1:D:264:GLU:C	2.62	0.41
2:E:32:PRO:HD2	2:E:84:HIS:CE1	2.55	0.41
1:D:66:LYS:HD3	3:F:4:MET:HA	2.03	0.41
2:E:37:VAL:HB	2:E:66:TYR:CZ	2.55	0.41
4:H:8:GLY:HA3	4:H:20:LEU:CD2	2.51	0.41
4:H:37:TRP:HD1	4:H:70:ILE:HD12	1.85	0.41
4:H:71:SER:O	4:H:79:PHE:HB2	2.21	0.41
1:A:42:SER:O	1:A:44:ARG:HG2	2.20	0.41
1:D:47:PRO:HB3	1:D:60:TRP:CH2	2.55	0.41
4:G:106:PRO:C	4:G:108:ASP:H	2.28	0.41
1:A:168:LEU:O	1:A:168:LEU:HD12	2.21	0.41
1:D:166:GLU:O	1:D:170:ARG:HG3	2.20	0.41
1:D:215:LEU:HD12	1:D:215:LEU:N	2.36	0.41
4:H:24:VAL:HG12	4:H:27:TYR:CD1	2.55	0.41
3:F:5:TRP:HE1	4:H:51:SER:CB	2.33	0.41
1:A:217:TRP:HE1	1:A:245:ALA:C	2.28	0.40
1:A:50:PRO:HA	1:A:53:GLU:HG3	2.01	0.40
1:A:117:ALA:HB2	2:B:60:TRP:CD2	2.56	0.40
4:H:20:LEU:HD12	4:H:81:LEU:HD23	2.03	0.40
1:A:202:ARG:HD2	1:A:244:TRP:CD2	2.55	0.40
1:D:202:ARG:HD2	1:D:204:TRP:CE2	2.56	0.40
2:B:7:ILE:HG21	2:B:93:VAL:CG2	2.50	0.40
1:D:181:ARG:NH1	1:D:209:TYR:HB2	2.37	0.40
2:B:33:SER:O	2:B:34:ASP:CB	2.69	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	270/276 (98%)	258 (96%)	12 (4%)	0	100	100
1	D	235/276 (85%)	224 (95%)	11 (5%)	0	100	100
2	B	95/100 (95%)	88 (93%)	7 (7%)	0	100	100
2	E	96/100 (96%)	91 (95%)	5 (5%)	0	100	100
3	C	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	F	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
4	G	105/147 (71%)	97 (92%)	8 (8%)	0	100	100
4	H	118/147 (80%)	117 (99%)	1 (1%)	0	100	100
All	All	933/1064 (88%)	887 (95%)	46 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	221/232 (95%)	221 (100%)	0	100	100
1	D	203/232 (88%)	203 (100%)	0	100	100
2	B	90/95 (95%)	90 (100%)	0	100	100
2	E	92/95 (97%)	92 (100%)	0	100	100
3	C	9/9 (100%)	9 (100%)	0	100	100
3	F	9/9 (100%)	9 (100%)	0	100	100
4	G	93/125 (74%)	93 (100%)	0	100	100
4	H	101/125 (81%)	101 (100%)	0	100	100
All	All	818/922 (89%)	818 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	87	GLN
1	A	191	HIS
2	B	13	HIS
2	B	24	ASN
1	D	86	ASN
1	D	93	HIS
4	G	77	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	274/276 (99%)	-0.35	5 (1%) 67 59	23, 41, 83, 96	0
1	D	247/276 (89%)	-0.19	7 (2%) 55 46	27, 48, 98, 118	0
2	B	99/100 (99%)	-0.50	1 (1%) 79 73	20, 40, 62, 76	0
2	E	98/100 (98%)	-0.53	1 (1%) 79 73	24, 45, 73, 83	0
3	C	9/9 (100%)	-0.57	0 100 100	33, 37, 38, 41	0
3	F	9/9 (100%)	-0.81	0 100 100	30, 38, 44, 45	0
4	G	111/147 (75%)	-0.28	1 (0%) 81 75	30, 48, 71, 87	0
4	H	120/147 (81%)	-0.39	0 100 100	31, 49, 82, 94	0
All	All	967/1064 (90%)	-0.35	15 (1%) 70 62	20, 45, 85, 118	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	196	ASP	4.7
2	E	97	ARG	3.9
1	A	275	GLU	3.5
1	D	209	TYR	3.2
1	D	199	ALA	3.1
4	G	72	VAL	2.8
1	D	197	HIS	2.6
1	A	194	VAL	2.6
2	B	98	ASP	2.6
1	A	18	GLY	2.5
1	D	270	LEU	2.4
1	D	210	PRO	2.3
1	D	250	PRO	2.3
1	D	275	GLU	2.3
1	A	227	ASP	2.2

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

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($\text{\AA}^2$)	Q<0.9
5	CL	A	301	1/1	0.97	0.13	47,47,47,47	0
5	CL	D	301	1/1	0.97	0.14	44,44,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.