



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:35 PM UTC

PDB ID : 2P1Y / pdb_00002p1y
Title : 1.B2.D9, a bispecific alpha/beta TCR
Authors : McBeth, C.; Pizarro, J.C.; Strong, R.K.
Deposited on : 2007-03-06
Resolution : 2.42 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	:	9.0.010 (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

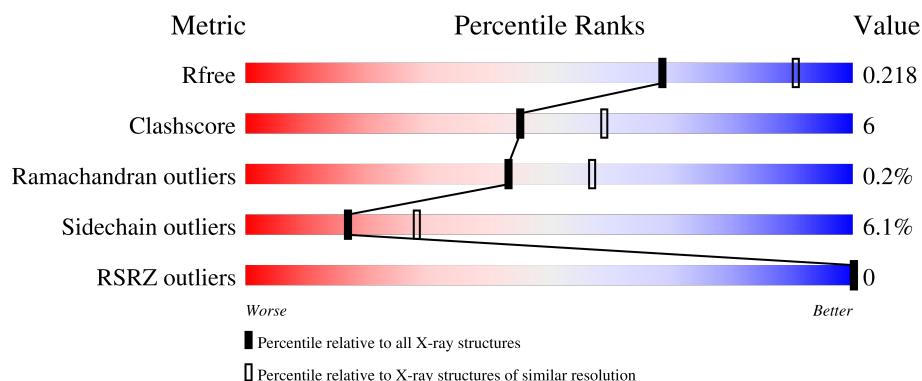
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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	238	 66% 27% . .
1	C	238	 62% 31% . .
1	E	238	 65% 22% . 9%
1	G	238	 67% 21% . 9%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7017 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 bispecific alpha/beta TCR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	231	Total	C	N	O	S	0	0	0
			1779	1124	304	346	5			
1	C	231	Total	C	N	O	S	0	0	0
			1771	1116	304	346	5			
1	E	217	Total	C	N	O	S	0	0	0
			1675	1064	283	323	5			
1	G	217	Total	C	N	O	S	0	0	0
			1690	1073	288	324	5			

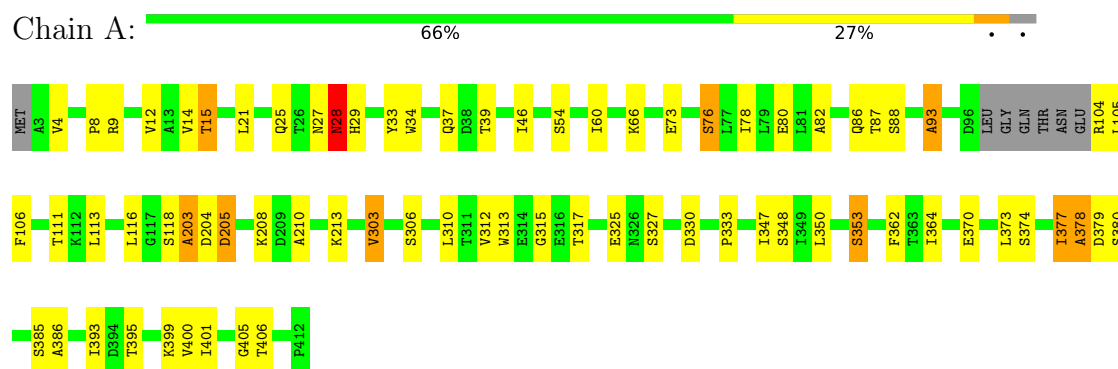
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	25	Total	O	0	0
			25	25		
2	C	15	Total	O	0	0
			15	15		
2	E	30	Total	O	0	0
			30	30		
2	G	32	Total	O	0	0
			32	32		

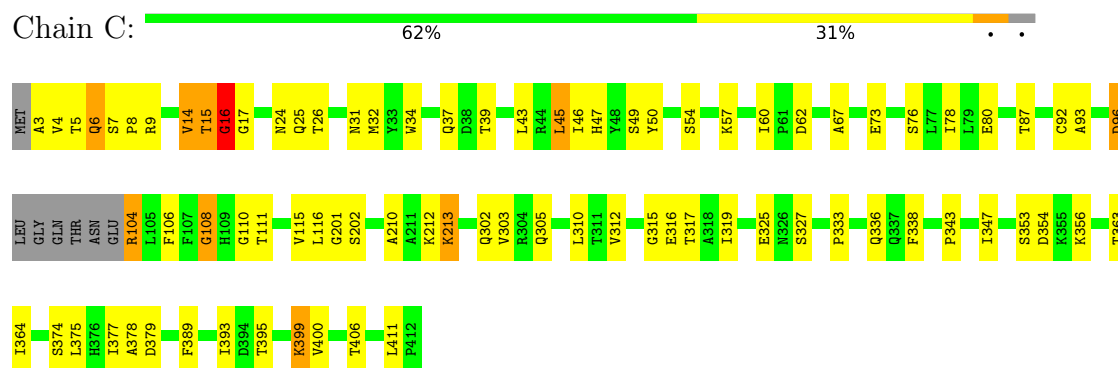
3 Residue-property plots

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.

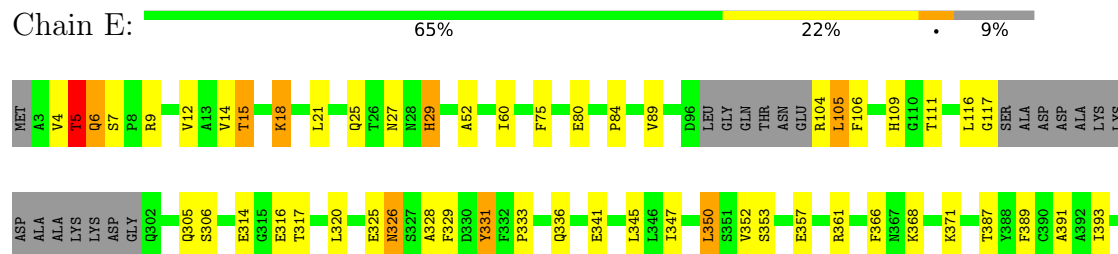
- Molecule 1: bispecific alpha/beta TCR



- Molecule 1: bispecific alpha/beta TCR

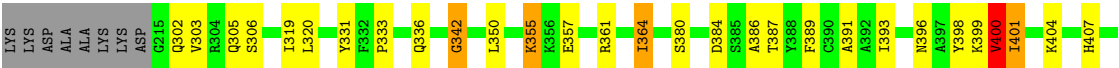
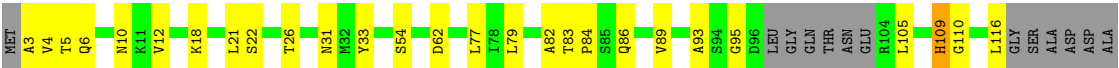


- Molecule 1: bispecific alpha/beta TCR





● Molecule 1: bispecific alpha/beta TCR



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.68Å 105.48Å 81.53Å 90.00° 116.04° 90.00°	Depositor
Resolution (Å)	40.39 – 2.42 40.39 – 2.42	Depositor EDS
% Data completeness (in resolution range)	99.4 (40.39-2.42) 99.2 (40.39-2.42)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.33 (at 2.42Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.181 , 0.254 0.188 , 0.218	Depositor DCC
R_{free} test set	2073 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	31.4	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 6.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.488 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7017	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.78% 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

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	2.01	31/1821 (1.7%)	1.52	10/2465 (0.4%)
1	C	2.07	43/1814 (2.4%)	1.61	22/2459 (0.9%)
1	E	1.82	28/1717 (1.6%)	1.52	13/2329 (0.6%)
1	G	1.75	19/1733 (1.1%)	1.48	15/2350 (0.6%)
All	All	1.92	121/7085 (1.7%)	1.53	60/9603 (0.6%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	2
1	E	0	2
1	G	0	1
All	All	0	6

All (121) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	4	VAL	CA-CB	11.35	1.68	1.54
1	C	4	VAL	CA-CB	10.37	1.67	1.54
1	E	4	VAL	CA-CB	10.16	1.67	1.53
1	C	78	ILE	CA-CB	9.66	1.66	1.54
1	A	364	ILE	C-O	-8.95	1.13	1.24
1	A	78	ILE	CA-CB	8.90	1.65	1.54
1	E	89	VAL	N-CA	8.10	1.56	1.46
1	C	45	LEU	C-O	7.73	1.33	1.24
1	C	336	GLN	N-CA	-7.63	1.36	1.46
1	E	345	LEU	C-O	7.31	1.32	1.23
1	G	4	VAL	CA-CB	7.25	1.63	1.53
1	C	60	ILE	CA-CB	7.22	1.63	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	377	ILE	CA-CB	7.19	1.62	1.53
1	G	386	ALA	CA-CB	7.19	1.67	1.53
1	C	73	GLU	CG-CD	6.89	1.69	1.52
1	G	355	LYS	CA-C	6.89	1.60	1.52
1	C	87	THR	CA-CB	6.87	1.64	1.53
1	G	95	GLY	C-O	-6.78	1.17	1.23
1	G	12	VAL	CA-CB	6.77	1.62	1.54
1	C	312	VAL	CA-CB	6.68	1.66	1.54
1	E	389	PHE	N-CA	6.68	1.54	1.46
1	G	336	GLN	N-CA	-6.62	1.38	1.46
1	C	62	ASP	C-O	6.58	1.31	1.23
1	A	205	ASP	CA-C	6.48	1.61	1.52
1	A	12	VAL	CA-CB	6.44	1.62	1.54
1	C	399	LYS	N-CA	6.42	1.54	1.46
1	C	377	ILE	CA-CB	6.41	1.61	1.53
1	C	9	ARG	CB-CG	6.37	1.71	1.52
1	C	315	GLY	C-O	-6.25	1.14	1.24
1	A	312	VAL	CA-CB	6.23	1.64	1.54
1	C	364	ILE	C-O	-6.21	1.16	1.24
1	E	29	HIS	C-O	-6.21	1.16	1.23
1	A	385	SER	C-O	6.20	1.31	1.24
1	A	60	ILE	CA-CB	6.17	1.62	1.54
1	A	203	ALA	CA-CB	6.16	1.63	1.53
1	E	387	THR	N-CA	6.16	1.53	1.46
1	C	108	GLY	C-O	-6.14	1.18	1.23
1	A	315	GLY	C-O	-6.14	1.15	1.24
1	C	375	LEU	CA-C	6.14	1.60	1.52
1	E	336	GLN	N-CA	-6.11	1.39	1.46
1	E	409	HIS	CA-C	6.08	1.60	1.52
1	A	210	ALA	CA-C	6.08	1.61	1.52
1	C	8	PRO	N-CA	6.07	1.57	1.47
1	G	387	THR	N-CA	6.06	1.53	1.46
1	E	341	GLU	C-O	6.05	1.31	1.23
1	C	347	ILE	C-O	-6.05	1.16	1.24
1	C	9	ARG	CD-NE	6.04	1.54	1.46
1	E	314	GLU	CA-C	6.02	1.60	1.52
1	E	409	HIS	C-O	-6.00	1.17	1.24
1	A	353	SER	N-CA	5.98	1.52	1.45
1	E	105	LEU	N-CA	5.97	1.53	1.46
1	C	6	GLN	C-O	-5.95	1.16	1.23
1	C	317	THR	CA-CB	5.95	1.61	1.53
1	C	343	PRO	C-O	5.95	1.30	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	111	THR	CA-CB	5.94	1.62	1.53
1	A	373	LEU	C-O	-5.92	1.16	1.23
1	A	370	GLU	C-O	-5.91	1.16	1.24
1	E	75	PHE	C-O	5.89	1.30	1.23
1	C	24	ASN	CA-C	-5.87	1.45	1.52
1	A	73	GLU	CG-CD	5.84	1.66	1.52
1	G	89	VAL	N-CA	5.82	1.53	1.46
1	G	93	ALA	CA-C	5.81	1.59	1.52
1	G	77	LEU	C-O	-5.79	1.17	1.24
1	A	9	ARG	CB-CG	5.77	1.69	1.52
1	C	406	THR	CA-C	-5.77	1.45	1.52
1	A	33	TYR	C-O	-5.72	1.17	1.23
1	A	37	GLN	C-O	-5.70	1.17	1.24
1	C	213	LYS	N-CA	5.69	1.53	1.46
1	E	398	TYR	CA-CB	5.66	1.62	1.53
1	A	93	ALA	CA-CB	-5.64	1.42	1.54
1	E	5	THR	C-O	-5.64	1.17	1.23
1	C	115	VAL	CA-CB	5.64	1.60	1.54
1	G	387	THR	CA-CB	-5.64	1.45	1.53
1	C	333	PRO	C-O	5.60	1.30	1.23
1	E	15	THR	CA-C	5.59	1.59	1.52
1	G	389	PHE	N-CA	5.58	1.52	1.46
1	E	12	VAL	CA-CB	5.57	1.62	1.54
1	G	33	TYR	N-CA	5.56	1.52	1.46
1	A	406	THR	CA-C	-5.55	1.46	1.52
1	E	410	VAL	C-O	-5.51	1.17	1.24
1	A	76	SER	CA-C	5.49	1.60	1.52
1	A	386	ALA	C-O	5.44	1.29	1.23
1	A	364	ILE	N-CA	5.43	1.53	1.46
1	C	305	GLN	C-O	5.42	1.30	1.23
1	C	37	GLN	N-CA	-5.39	1.39	1.46
1	C	34	TRP	C-O	-5.38	1.17	1.24
1	G	21	LEU	C-O	5.36	1.30	1.24
1	G	364	ILE	N-CA	5.34	1.52	1.46
1	E	405	GLY	N-CA	5.34	1.51	1.45
1	C	347	ILE	CB-CG2	5.33	1.70	1.52
1	C	389	PHE	C-O	-5.33	1.17	1.24
1	C	353	SER	C-O	-5.33	1.17	1.23
1	C	319	ILE	CA-CB	5.32	1.60	1.54
1	C	50	TYR	CA-CB	5.31	1.62	1.53
1	E	52	ALA	C-O	5.31	1.30	1.23
1	A	313	TRP	N-CA	5.29	1.52	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	331	TYR	CA-C	5.24	1.59	1.52
1	C	7	SER	CA-C	5.21	1.59	1.52
1	E	60	ILE	CA-CB	5.21	1.60	1.54
1	G	10	ASN	N-CA	5.17	1.52	1.46
1	C	37	GLN	C-O	-5.15	1.18	1.24
1	A	34	TRP	C-O	-5.15	1.18	1.24
1	A	378	ALA	CA-CB	5.14	1.61	1.53
1	E	387	THR	CA-CB	-5.14	1.46	1.53
1	E	398	TYR	C-N	5.14	1.40	1.33
1	C	363	THR	N-CA	5.11	1.52	1.46
1	E	6	GLN	C-O	-5.10	1.17	1.23
1	A	317	THR	CA-CB	5.08	1.59	1.53
1	C	202	SER	N-CA	5.08	1.52	1.45
1	A	113	LEU	CA-C	5.06	1.58	1.52
1	A	46	ILE	CA-C	5.05	1.59	1.52
1	C	354	ASP	CA-C	5.05	1.59	1.52
1	E	18	LYS	CA-C	5.05	1.58	1.52
1	A	380	SER	CA-C	5.04	1.59	1.52
1	C	210	ALA	N-CA	5.04	1.52	1.46
1	G	82	ALA	CA-CB	5.03	1.61	1.53
1	G	319	ILE	C-O	-5.03	1.18	1.24
1	E	60	ILE	CA-C	5.02	1.58	1.52
1	G	364	ILE	CA-C	5.02	1.58	1.52
1	C	16	GLY	CA-C	5.02	1.57	1.52
1	E	361	ARG	N-CA	5.01	1.52	1.46

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	15	THR	CA-C-N	-12.36	106.19	120.03
1	C	15	THR	C-N-CA	-12.36	106.19	120.03
1	G	411	LEU	N-CA-C	11.13	126.23	110.20
1	E	411	LEU	N-CA-C	10.91	125.92	110.20
1	G	306	SER	N-CA-C	10.90	125.99	110.40
1	E	105	LEU	N-CA-C	8.81	123.67	110.64
1	C	6	GLN	N-CA-C	8.76	123.48	110.52
1	G	302	GLN	N-CA-C	8.67	123.32	109.79
1	G	364	ILE	N-CA-C	7.89	119.26	108.84
1	E	306	SER	N-CA-C	7.83	121.60	110.40
1	C	202	SER	N-CA-C	7.83	119.90	110.19
1	C	338	PHE	CA-C-N	7.62	127.61	119.76
1	C	338	PHE	C-N-CA	7.62	127.61	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	306	SER	N-CA-CB	-7.38	97.91	111.19
1	E	316	GLU	N-CA-C	-7.35	100.83	110.53
1	C	16	GLY	N-CA-C	7.35	121.79	112.83
1	A	405	GLY	N-CA-C	7.33	121.34	112.48
1	A	306	SER	N-CA-C	7.25	121.07	108.82
1	A	303	VAL	N-CA-C	-6.90	98.45	108.11
1	E	5	THR	CB-CA-C	6.87	121.03	109.48
1	C	14	VAL	CB-CA-C	6.74	121.05	110.81
1	C	5	THR	CB-CA-C	6.71	120.87	110.14
1	E	6	GLN	N-CA-C	6.70	120.38	110.46
1	C	39	THR	CA-C-N	-6.54	112.92	122.46
1	C	39	THR	C-N-CA	-6.54	112.92	122.46
1	A	364	ILE	N-CA-C	6.37	117.75	108.58
1	G	380	SER	N-CA-C	6.20	118.31	110.24
1	E	400	VAL	CB-CA-C	-6.19	102.07	110.42
1	C	364	ILE	N-CA-C	6.12	117.39	108.58
1	E	5	THR	N-CA-C	-6.02	99.89	109.76
1	E	397	ALA	CA-C-N	5.88	132.47	123.47
1	E	397	ALA	C-N-CA	5.88	132.47	123.47
1	G	62	ASP	N-CA-C	5.72	118.52	110.23
1	C	17	GLY	N-CA-C	5.72	121.30	111.98
1	G	400	VAL	CB-CA-C	-5.70	101.94	111.29
1	G	5	THR	N-CA-C	-5.67	100.47	109.59
1	A	347	ILE	CA-C-O	-5.64	115.66	120.96
1	C	303	VAL	N-CA-C	-5.63	100.29	108.17
1	A	28	ASN	N-CA-CB	-5.62	103.92	112.47
1	C	316	GLU	CA-CB-CG	5.62	125.33	114.10
1	A	105	LEU	CA-C-N	5.61	130.89	122.99
1	A	105	LEU	C-N-CA	5.61	130.89	122.99
1	G	361	ARG	NE-CZ-NH2	5.56	124.21	119.20
1	G	109	HIS	CB-CA-C	5.49	119.84	110.56
1	C	57	LYS	N-CA-C	-5.41	102.27	110.28
1	C	9	ARG	NE-CZ-NH2	5.35	124.02	119.20
1	E	326	ASN	CB-CA-C	5.34	119.32	110.78
1	G	342	GLY	CA-C-N	-5.30	114.32	119.78
1	G	342	GLY	C-N-CA	-5.30	114.32	119.78
1	E	317	THR	N-CA-C	-5.27	100.58	108.96
1	A	4	VAL	N-CA-C	-5.27	100.73	108.11
1	G	384	ASP	N-CA-C	-5.16	106.93	113.18
1	E	347	ILE	N-CA-C	-5.16	100.77	107.88
1	G	54	SER	N-CA-C	5.13	116.78	109.14
1	G	303	VAL	CB-CA-C	5.12	116.25	110.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	201	GLY	N-CA-C	-5.08	101.15	113.18
1	C	110	GLY	N-CA-C	5.05	117.94	112.29
1	C	15	THR	O-C-N	-5.04	115.89	122.59
1	C	43	LEU	CA-C-N	-5.01	115.43	122.94
1	C	43	LEU	C-N-CA	-5.01	115.43	122.94

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	104	ARG	Peptide
1	C	116	LEU	Peptide
1	C	16	GLY	Peptide
1	E	305	GLN	Peptide
1	E	5	THR	Peptide
1	G	305	GLN	Peptide

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	1779	0	1697	24	0
1	C	1771	0	1671	18	0
1	E	1675	0	1590	23	0
1	G	1690	0	1608	15	0
2	A	25	0	0	0	0
2	C	15	0	0	1	0
2	E	30	0	0	0	0
2	G	32	0	0	1	0
All	All	7017	0	6566	79	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 6.

All (79) 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:104:ARG:O	1:C:104:ARG:HG3	1.64	0.98
1:E:25:GLN:HE22	1:E:29:HIS:H	1.12	0.94
1:C:16:GLY:HA3	1:G:3:ALA:N	1.88	0.88
1:A:15:THR:HG22	1:A:116:LEU:O	1.79	0.82
1:C:96:ASP:OD1	1:C:96:ASP:C	2.25	0.80
1:G:109:HIS:CD2	1:G:342:GLY:H	2.02	0.77
1:A:330:ASP:OD1	1:A:395:THR:HG23	1.86	0.75
1:A:303:VAL:HG23	1:A:401:ILE:HG22	1.68	0.73
1:G:109:HIS:HD2	1:G:342:GLY:H	1.39	0.67
1:A:213:LYS:HG2	1:A:325:GLU:HG3	1.75	0.67
1:G:407:HIS:CE1	2:G:437:HOH:O	2.48	0.67
1:C:3:ALA:HA	1:C:26:THR:OG1	1.97	0.64
1:A:118:SER:C	1:A:203:ALA:N	2.56	0.64
1:G:333:PRO:HG2	1:G:391:ALA:HB3	1.79	0.63
1:C:104:ARG:O	1:C:104:ARG:CG	2.44	0.63
1:E:5:THR:HA	1:E:109:HIS:HE1	1.65	0.61
1:E:329:PHE:CZ	1:E:401:ILE:HG13	2.36	0.61
1:C:93:ALA:HA	1:C:106:PHE:O	2.00	0.60
1:A:303:VAL:HG23	1:A:401:ILE:CG2	2.31	0.60
1:A:66:LYS:HD2	1:A:80:GLU:OE2	2.00	0.59
1:E:333:PRO:HG2	1:E:391:ALA:HB3	1.84	0.59
1:A:27:ASN:O	1:A:28:ASN:CB	2.51	0.58
1:G:83:THR:OG1	1:G:86:GLN:HG3	2.03	0.58
1:A:303:VAL:CG2	1:A:401:ILE:HG22	2.34	0.58
1:E:9:ARG:CD	1:E:109:HIS:CD2	2.86	0.57
1:E:326:ASN:ND2	1:E:328:ALA:H	2.04	0.55
1:A:362:PHE:CE2	1:A:377:ILE:HG12	2.42	0.55
1:E:15:THR:OG1	1:E:117:GLY:HA3	2.07	0.55
1:C:6:GLN:HE21	1:C:108:GLY:HA3	1.73	0.54
1:G:79:LEU:HD23	1:G:86:GLN:OE1	2.08	0.54
1:A:93:ALA:HA	1:A:106:PHE:O	2.08	0.53
1:C:212:LYS:O	1:C:302:GLN:HG3	2.08	0.53
1:E:6:GLN:H	1:E:109:HIS:CE1	2.27	0.53
1:C:15:THR:O	1:C:16:GLY:C	2.47	0.52
1:G:26:THR:HG22	1:G:26:THR:O	2.11	0.51
1:C:6:GLN:NE2	1:C:92:CYS:H	2.09	0.51
1:C:80:GLU:HG2	2:C:417:HOH:O	2.11	0.50
1:C:25:GLN:HG2	1:C:32:MET:HE3	1.94	0.49
1:A:204:ASP:O	1:A:208:LYS:HG3	2.12	0.49
1:C:213:LYS:HD3	1:C:325:GLU:HG3	1.95	0.48
1:E:331:TYR:HD2	1:E:393:ILE:HB	1.77	0.48
1:C:31:ASN:HA	1:C:49:SER:O	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:393:ILE:HG12	1:C:400:VAL:HG22	1.95	0.48
1:E:18:LYS:HD2	1:E:80:GLU:HA	1.95	0.48
1:E:21:LEU:HD22	1:E:111:THR:HG21	1.95	0.48
1:G:31:ASN:HD21	1:G:398:TYR:HE1	1.63	0.47
1:A:8:PRO:O	1:A:111:THR:HG23	2.15	0.47
1:E:325:GLU:O	1:E:371:LYS:NZ	2.41	0.46
1:C:46:ILE:HG22	1:C:47:HIS:CD2	2.51	0.46
1:A:39:THR:HG23	1:A:39:THR:O	2.14	0.46
1:A:118:SER:HB2	1:A:205:ASP:H	1.79	0.46
1:A:333:PRO:HA	1:A:348:SER:HB3	1.97	0.46
1:E:25:GLN:NE2	1:E:29:HIS:H	1.96	0.46
1:G:411:LEU:HB3	1:G:412:PRO:HA	1.97	0.45
1:E:391:ALA:HA	1:E:401:ILE:O	2.16	0.45
1:A:378:ALA:O	1:A:379:ASP:C	2.59	0.44
1:E:329:PHE:HZ	1:E:401:ILE:HG13	1.80	0.44
1:E:14:VAL:HA	1:E:116:LEU:O	2.18	0.44
1:G:400:VAL:C	1:G:401:ILE:HD13	2.42	0.43
1:A:39:THR:O	1:A:39:THR:CG2	2.67	0.43
1:E:25:GLN:HE21	1:E:27:ASN:H	1.67	0.43
1:A:87:THR:O	1:A:88:SER:HB2	2.18	0.43
1:G:6:GLN:HE22	1:G:110:GLY:HA2	1.83	0.43
1:G:331:TYR:HD2	1:G:393:ILE:HB	1.84	0.43
1:A:27:ASN:HB3	1:A:29:HIS:CE1	2.54	0.42
1:E:366:PHE:HE2	1:E:368:LYS:HB2	1.83	0.42
1:A:21:LEU:HD22	1:A:111:THR:HG21	2.00	0.42
1:E:25:GLN:HE21	1:E:27:ASN:HB2	1.83	0.42
1:C:378:ALA:O	1:C:379:ASP:C	2.60	0.42
1:E:25:GLN:NE2	1:E:29:HIS:HB2	2.35	0.41
1:G:355:LYS:HA	1:G:364:ILE:O	2.20	0.41
1:A:82:ALA:HA	1:A:86:GLN:OE1	2.21	0.41
1:A:400:VAL:O	1:A:401:ILE:HD13	2.21	0.41
1:E:104:ARG:O	1:E:106:PHE:CD2	2.74	0.41
1:A:362:PHE:HE2	1:A:377:ILE:HG12	1.86	0.40
1:E:350:LEU:C	1:E:352:VAL:H	2.29	0.40
1:E:350:LEU:C	1:E:352:VAL:N	2.77	0.40
1:G:6:GLN:NE2	1:G:110:GLY:C	2.80	0.40
1:C:67:ALA:HA	1:C:76:SER:O	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	225/238 (94%)	216 (96%)	9 (4%)	0	100	100
1	C	227/238 (95%)	216 (95%)	11 (5%)	0	100	100
1	E	211/238 (89%)	202 (96%)	8 (4%)	1 (0%)	24	35
1	G	211/238 (89%)	198 (94%)	12 (6%)	1 (0%)	24	35
All	All	874/952 (92%)	832 (95%)	40 (5%)	2 (0%)	43	57

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	84	PRO
1	E	84	PRO

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	190/199 (96%)	177 (93%)	13 (7%)	14	24
1	C	188/199 (94%)	176 (94%)	12 (6%)	16	26
1	E	180/199 (90%)	172 (96%)	8 (4%)	25	41
1	G	182/199 (92%)	170 (93%)	12 (7%)	15	25
All	All	740/796 (93%)	695 (94%)	45 (6%)	17	28

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	VAL
1	A	15	THR
1	A	25	GLN
1	A	28	ASN
1	A	54	SER
1	A	76	SER
1	A	310	LEU
1	A	327	SER
1	A	350	LEU
1	A	353	SER
1	A	374	SER
1	A	393	ILE
1	A	399	LYS
1	C	14	VAL
1	C	45	LEU
1	C	54	SER
1	C	96	ASP
1	C	104	ARG
1	C	310	LEU
1	C	327	SER
1	C	356	LYS
1	C	374	SER
1	C	395	THR
1	C	399	LYS
1	C	411	LEU
1	E	7	SER
1	E	105	LEU
1	E	320	LEU
1	E	350	LEU
1	E	353	SER
1	E	357	GLU
1	E	399	LYS
1	E	401	ILE
1	G	18	LYS
1	G	22	SER
1	G	105	LEU
1	G	116	LEU
1	G	320	LEU
1	G	350	LEU
1	G	357	GLU
1	G	396	ASN
1	G	399	LYS
1	G	400	VAL

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Mol	Chain	Res	Type
1	G	401	ILE
1	G	404	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	HIS
1	A	409	HIS
1	C	6	GLN
1	C	28	ASN
1	C	30	ASN
1	C	72	GLN
1	C	302	GLN
1	C	407	HIS
1	E	24	ASN
1	E	25	GLN
1	E	28	ASN
1	E	74	ASN
1	E	109	HIS
1	E	326	ASN
1	E	336	GLN
1	E	396	ASN
1	G	28	ASN
1	G	41	HIS
1	G	109	HIS
1	G	376	HIS
1	G	407	HIS

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 [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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	118:SER	C	203:ALA	N	2.56

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	231/238 (97%)	-1.51	0 100 100	11, 23, 54, 67	0
1	C	231/238 (97%)	-1.51	0 100 100	11, 23, 55, 64	0
1	E	217/238 (91%)	-1.51	0 100 100	12, 28, 50, 59	0
1	G	217/238 (91%)	-1.46	0 100 100	12, 28, 51, 58	0
All	All	896/952 (94%)	-1.50	0 100 100	11, 26, 53, 67	0

There are no RSRZ outliers to report.

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.