



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:08 PM UTC

PDB ID : 3R06 / pdb_00003r06
Title : Crystal structure of anti-mouse CD3epsilon antibody 2C11 Fab fragment
Authors : Shore, D.A.; Zhu, X.; Wilson, I.A.
Deposited on : 2011-03-07
Resolution : 2.50 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

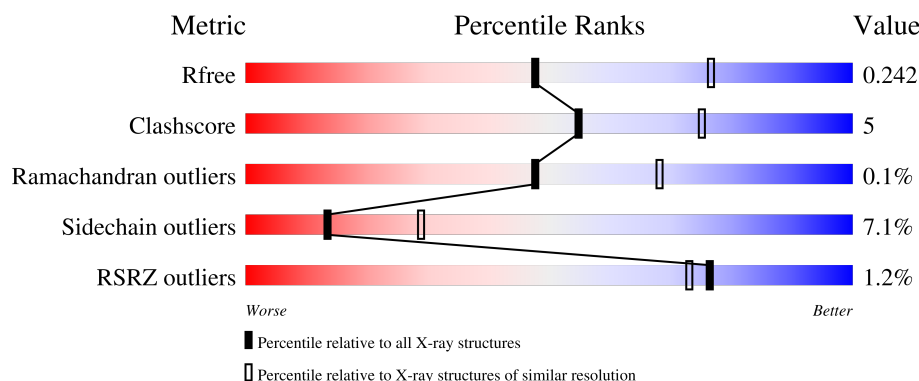
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	213	85% 14% .
1	C	213	86% 13% .
1	E	213	80% 18% .
1	L	213	88% 12% .
2	B	216	2% 80% 15% .

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Mol	Chain	Length	Quality of chain
2	D	216	% 81% 14% • •
2	F	216	% 75% 19% • •
2	H	216	3% 77% 17% • •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13427 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 anti-mouse CD3epsilon antibody 2C11 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	213	Total	C	N	O	S	0	0	0
			1664	1041	278	340	5			
1	C	213	Total	C	N	O	S	0	0	0
			1664	1041	278	340	5			
1	E	213	Total	C	N	O	S	0	0	0
			1664	1041	278	340	5			
1	L	213	Total	C	N	O	S	0	0	0
			1664	1041	278	340	5			

- Molecule 2 is a protein called anti-mouse CD3epsilon antibody 2C11 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	210	Total	C	N	O	S	0	0	0
			1586	1008	262	308	8			
2	D	211	Total	C	N	O	S	0	0	0
			1592	1011	263	309	9			
2	F	211	Total	C	N	O	S	0	0	0
			1592	1011	263	309	9			
2	H	211	Total	C	N	O	S	0	0	0
			1592	1011	263	309	9			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	56	Total	O	0	0
			56	56		
3	B	40	Total	O	0	0
			40	40		
3	C	63	Total	O	0	0
			63	63		
3	D	58	Total	O	0	0
			58	58		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	46	Total 46	O 46	0	0
3	F	39	Total 39	O 39	0	0
3	H	44	Total 44	O 44	0	0
3	L	63	Total 63	O 63	0	0

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.

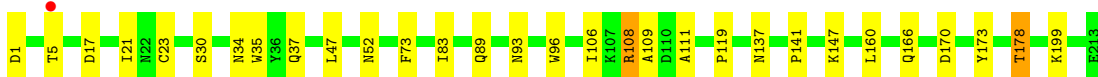
- Molecule 1: anti-mouse CD3epsilon antibody 2C11 Fab light chain

Chain A: 



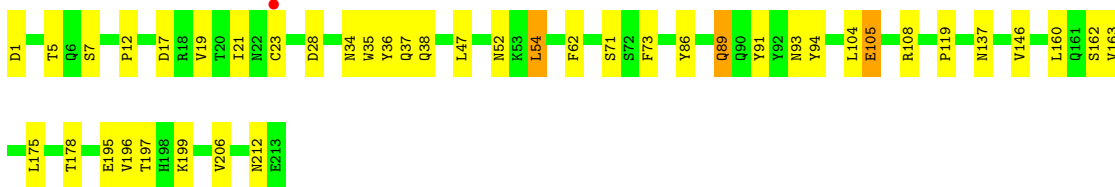
- Molecule 1: anti-mouse CD3epsilon antibody 2C11 Fab light chain

Chain C: 



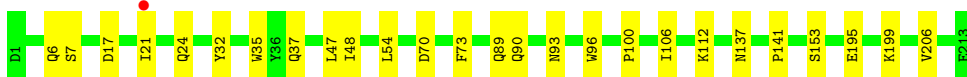
- Molecule 1: anti-mouse CD3epsilon antibody 2C11 Fab light chain

Chain E: 



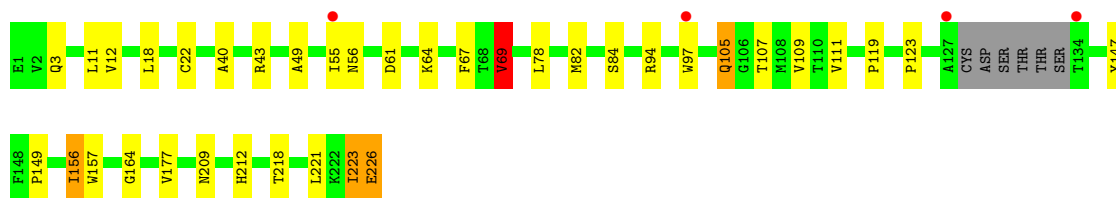
- Molecule 1: anti-mouse CD3epsilon antibody 2C11 Fab light chain

Chain L: 

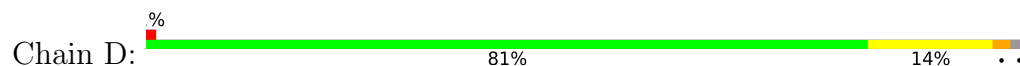


- Molecule 2: anti-mouse CD3epsilon antibody 2C11 Fab heavy chain

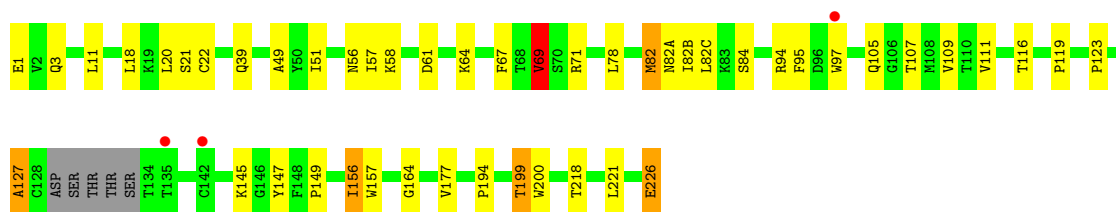
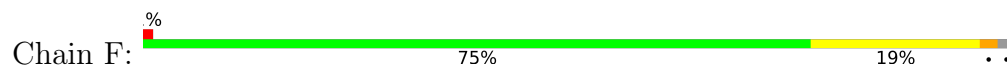
Chain B: 



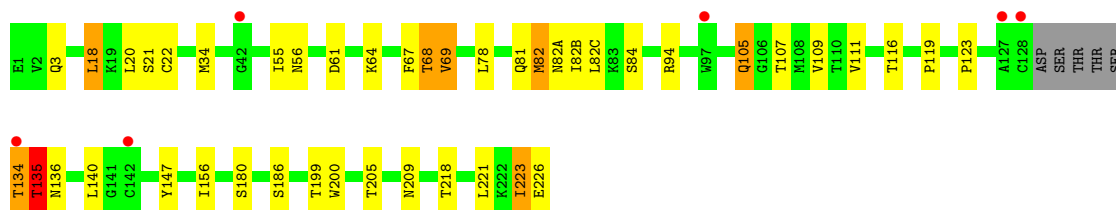
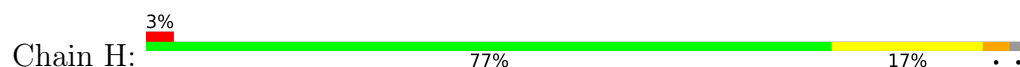
- Molecule 2: anti-mouse CD3epsilon antibody 2C11 Fab heavy chain



- Molecule 2: anti-mouse CD3epsilon antibody 2C11 Fab heavy chain



- Molecule 2: anti-mouse CD3epsilon antibody 2C11 Fab heavy chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	106.17Å 72.39Å 127.49Å 90.00° 109.52° 90.00°	Depositor
Resolution (Å)	36.02 – 2.50 36.02 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (36.02-2.50) 97.8 (36.02-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 2.51Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, R_{free}	0.191 , 0.236 0.199 , 0.242	Depositor DCC
R_{free} test set	3142 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	49.6	Xtriage
Anisotropy	0.469	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 64.6	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	0.95	EDS
Total number of atoms	13427	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.58% 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	0.82	0/1701	1.19	7/2309 (0.3%)
1	C	0.84	0/1701	1.17	7/2309 (0.3%)
1	E	0.80	0/1701	1.20	3/2309 (0.1%)
1	L	0.83	0/1701	1.21	7/2309 (0.3%)
2	B	0.77	0/1625	1.15	4/2216 (0.2%)
2	D	0.84	0/1631	1.22	7/2224 (0.3%)
2	F	0.77	0/1631	1.16	5/2224 (0.2%)
2	H	0.78	1/1631 (0.1%)	1.18	6/2224 (0.3%)
All	All	0.81	1/13322 (0.0%)	1.19	46/18124 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	34	MET	SD-CE	6.54	1.96	1.79

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	223	ILE	CA-C-N	7.00	134.30	121.70
2	H	223	ILE	C-N-CA	7.00	134.30	121.70
1	C	137	ASN	N-CA-C	6.89	120.90	109.95
1	L	106	ILE	N-CA-C	6.70	118.48	108.23
1	A	17	ASP	CA-CB-CG	6.63	119.23	112.60
2	H	134	THR	CA-C-N	6.61	134.16	121.54
2	H	134	THR	C-N-CA	6.61	134.16	121.54
1	E	17	ASP	CA-CB-CG	6.59	119.19	112.60
1	E	137	ASN	N-CA-C	6.44	120.55	110.17
1	A	137	ASN	N-CA-C	6.42	120.16	109.95
1	L	137	ASN	N-CA-C	6.40	120.47	110.17
2	D	127	ALA	CA-C-N	6.27	132.99	121.70
2	D	127	ALA	C-N-CA	6.27	132.99	121.70
1	C	17	ASP	CA-CB-CG	6.07	118.67	112.60
1	L	17	ASP	CA-CB-CG	5.93	118.53	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	170	ASP	N-CA-C	5.92	120.20	112.92
1	L	70	ASP	N-CA-C	5.85	118.06	108.52
2	F	127	ALA	CA-C-N	5.80	132.13	121.70
2	F	127	ALA	C-N-CA	5.80	132.13	121.70
2	D	67	PHE	CA-CB-CG	5.79	119.59	113.80
2	D	204	GLN	CA-C-N	5.71	128.96	120.90
2	D	204	GLN	C-N-CA	5.71	128.96	120.90
2	H	67	PHE	CA-CB-CG	5.67	119.47	113.80
1	L	141	PRO	CA-C-N	5.58	128.32	120.28
1	L	141	PRO	C-N-CA	5.58	128.32	120.28
1	C	34	ASN	CA-CB-CG	5.56	118.16	112.60
2	F	67	PHE	CA-CB-CG	5.51	119.31	113.80
2	B	67	PHE	CA-CB-CG	5.34	119.14	113.80
1	L	70	ASP	CA-CB-CG	5.31	117.91	112.60
1	A	28	ASP	CA-CB-CG	5.30	117.91	112.60
2	B	223	ILE	CA-C-N	5.30	131.23	121.70
2	B	223	ILE	C-N-CA	5.30	131.23	121.70
2	F	69	VAL	N-CA-CB	5.29	117.86	110.56
1	A	108	ARG	N-CA-CB	-5.28	103.34	111.46
1	C	141	PRO	CA-C-N	5.24	127.82	120.28
1	C	141	PRO	C-N-CA	5.24	127.82	120.28
2	H	69	VAL	N-CA-CB	5.22	117.42	110.26
2	B	69	VAL	N-CA-CB	5.15	117.67	110.56
1	E	108	ARG	N-CA-CB	-5.12	103.58	111.46
2	D	69	VAL	N-CA-CB	5.11	117.60	110.56
1	A	141	PRO	CA-C-N	5.10	127.63	120.28
1	A	141	PRO	C-N-CA	5.10	127.63	120.28
1	C	147	LYS	N-CA-C	5.10	117.42	109.52
2	F	145	LYS	N-CA-C	5.06	118.06	109.76
1	A	137	ASN	CA-CB-CG	5.04	117.64	112.60
2	D	82(C)	LEU	N-CA-C	5.04	117.51	109.96

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	1664	0	1608	14	0
1	C	1664	0	1608	11	0
1	E	1664	0	1608	24	0
1	L	1664	0	1608	8	0
2	B	1586	0	1559	19	0
2	D	1592	0	1564	18	0
2	F	1592	0	1564	25	0
2	H	1592	0	1564	17	0
3	A	56	0	0	0	0
3	B	40	0	0	1	0
3	C	63	0	0	0	0
3	D	58	0	0	0	0
3	E	46	0	0	1	0
3	F	39	0	0	0	0
3	H	44	0	0	0	0
3	L	63	0	0	0	0
All	All	13427	0	12683	123	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 (123) 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)
2:D:51:ILE:HG22	2:D:57:ILE:HG12	1.43	1.00
2:F:51:ILE:HG22	2:F:57:ILE:HG12	1.42	0.99
2:F:51:ILE:HD12	2:F:71:ARG:HD2	1.55	0.89
2:D:51:ILE:HD12	2:D:71:ARG:HD2	1.57	0.87
1:E:38:GLN:HE22	2:F:39:GLN:HE22	1.25	0.83
2:H:156:ILE:HG13	2:H:209:ASN:HB2	1.64	0.80
2:H:200:TRP:CZ2	2:H:226:GLU:HB2	2.27	0.70
2:D:226:GLU:HG3	2:D:226:GLU:OXT	1.94	0.66
2:F:156:ILE:HG23	2:F:157:TRP:N	2.13	0.63
2:B:156:ILE:HG23	2:B:157:TRP:N	2.13	0.63
1:C:108:ARG:NH1	1:C:109:ALA:O	2.32	0.63
2:D:156:ILE:HG13	2:D:209:ASN:HB2	1.81	0.62
2:F:116:THR:HG22	2:F:149:PRO:HD3	1.82	0.62
1:A:37:GLN:HB2	1:A:47:LEU:HD11	1.83	0.61
1:L:89:GLN:HE21	1:L:96:TRP:HB3	1.67	0.60
2:B:82:MET:HE1	2:B:109:VAL:HG11	1.83	0.60
1:C:108:ARG:HH12	1:C:111:ALA:HB2	1.67	0.60
2:F:156:ILE:HG12	2:F:164:GLY:HA2	1.82	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:156:ILE:HG12	2:B:164:GLY:HA2	1.84	0.59
2:H:82:MET:HE1	2:H:109:VAL:HG11	1.83	0.59
1:E:19:VAL:HG11	1:E:104:LEU:HD21	1.85	0.59
2:H:134:THR:HG23	2:H:136:ASN:H	1.67	0.59
2:H:123:PRO:HD3	2:H:221:LEU:HD21	1.84	0.59
2:H:200:TRP:HZ2	2:H:226:GLU:HB2	1.66	0.59
2:D:51:ILE:HD11	2:D:78:LEU:CD1	2.33	0.58
1:E:36:TYR:OH	2:F:95:PHE:HE2	1.86	0.58
1:A:186:TYR:CZ	1:A:211:ARG:HG3	2.38	0.58
2:D:226:GLU:OXT	2:D:226:GLU:CG	2.51	0.57
2:F:51:ILE:HD11	2:F:78:LEU:CD1	2.34	0.57
1:C:89:GLN:HE21	1:C:96:TRP:HB3	1.70	0.57
1:E:37:GLN:HB2	1:E:47:LEU:HD11	1.87	0.56
2:D:82:MET:HE1	2:D:109:VAL:HG21	1.86	0.56
2:F:119:PRO:HB3	2:F:147:TYR:HB3	1.88	0.55
2:B:119:PRO:HB3	2:B:147:TYR:HB3	1.88	0.55
2:B:149:PRO:O	2:B:212:HIS:HE1	1.91	0.54
2:H:156:ILE:CG1	2:H:209:ASN:HB2	2.36	0.54
1:A:36:TYR:HE1	1:A:89:GLN:HG2	1.72	0.53
1:E:54:LEU:HD11	1:E:62:PHE:O	2.08	0.53
1:E:36:TYR:HE1	1:E:89:GLN:HG2	1.73	0.53
1:C:166:GLN:HG3	1:C:173:TYR:CZ	2.45	0.52
1:E:38:GLN:NE2	2:F:39:GLN:HE22	2.03	0.52
1:A:160:LEU:HD22	2:B:177:VAL:HG11	1.92	0.51
2:B:223:ILE:O	2:B:226:GLU:HB2	2.10	0.51
1:E:34:ASN:ND2	1:E:91:TYR:CE1	2.78	0.51
2:B:156:ILE:HG12	3:B:316:HOH:O	2.11	0.50
2:D:156:ILE:CG1	2:D:209:ASN:HB2	2.41	0.50
2:F:123:PRO:HD3	2:F:221:LEU:HD21	1.93	0.50
1:A:160:LEU:HD22	2:B:177:VAL:CG1	2.43	0.49
2:D:140:LEU:HB3	2:D:223:ILE:HG21	1.94	0.49
1:A:34:ASN:HD22	1:A:89:GLN:HE21	1.59	0.49
1:C:37:GLN:HB2	1:C:47:LEU:HD11	1.95	0.48
1:E:38:GLN:HE22	2:F:39:GLN:NE2	2.04	0.48
2:H:134:THR:HG23	2:H:135:THR:H	1.78	0.48
2:B:157:TRP:HZ3	2:B:223:ILE:HD13	1.78	0.48
2:H:18:LEU:HB3	2:H:82:MET:HE2	1.95	0.48
1:L:37:GLN:HB2	1:L:47:LEU:HD11	1.95	0.48
2:B:123:PRO:HD3	2:B:221:LEU:HD21	1.96	0.48
1:A:195:GLU:HG3	1:A:206:VAL:HG22	1.96	0.47
1:E:119:PRO:HG3	2:F:127:ALA:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:160:LEU:HD22	2:F:177:VAL:CG1	2.44	0.47
1:C:119:PRO:HG3	2:D:127:ALA:HB2	1.97	0.47
2:D:123:PRO:HD3	2:D:221:LEU:HD21	1.96	0.47
1:E:89:GLN:O	1:E:89:GLN:HG3	2.15	0.46
1:E:160:LEU:HD22	2:F:177:VAL:HG11	1.95	0.46
2:F:200:TRP:HZ2	2:F:226:GLU:HB2	1.80	0.46
2:H:140:LEU:HB3	2:H:223:ILE:HG21	1.98	0.46
2:F:61:ASP:HA	2:F:64:LYS:HD2	1.98	0.46
2:B:156:ILE:HG22	2:B:209:ASN:OD1	2.16	0.45
1:C:160:LEU:HB2	1:C:178:THR:HG22	1.98	0.45
2:H:119:PRO:HB3	2:H:147:TYR:HB3	1.98	0.45
1:E:160:LEU:HB3	2:F:177:VAL:HG11	1.98	0.45
1:L:35:TRP:CE2	1:L:73:PHE:HB2	2.52	0.45
2:F:49:ALA:HB1	2:F:69:VAL:HG21	1.99	0.45
2:H:22:CYS:HB3	2:H:78:LEU:HB3	1.98	0.45
2:D:61:ASP:HA	2:D:64:LYS:HD2	1.99	0.44
1:E:162:SER:O	1:E:175:LEU:HD12	2.17	0.44
1:E:178:THR:HG23	3:E:301:HOH:O	2.16	0.44
2:H:61:ASP:HA	2:H:64:LYS:HD2	1.98	0.44
2:H:68:THR:HG23	2:H:81:GLN:HB3	2.00	0.44
1:L:195:GLU:HG3	1:L:206:VAL:HG22	1.99	0.44
2:B:49:ALA:HB1	2:B:69:VAL:HG21	1.99	0.44
1:A:89:GLN:O	1:A:89:GLN:HG3	2.17	0.43
2:B:61:ASP:HA	2:B:64:LYS:HD2	1.98	0.43
2:D:22:CYS:HB3	2:D:78:LEU:HB3	1.99	0.43
1:E:86:TYR:CE1	1:E:104:LEU:HD12	2.53	0.43
1:C:23:CYS:HB2	1:C:35:TRP:CH2	2.53	0.43
2:H:105:GLN:H	2:H:105:GLN:HG3	1.53	0.43
2:B:84:SER:HA	2:B:111:VAL:HB	2.01	0.43
2:F:22:CYS:HB3	2:F:78:LEU:HB3	2.01	0.43
2:H:84:SER:HA	2:H:111:VAL:HB	2.01	0.43
2:H:123:PRO:CD	2:H:221:LEU:HD21	2.48	0.43
2:D:119:PRO:HB3	2:D:147:TYR:HB3	2.01	0.42
2:F:82:MET:HE1	2:F:109:VAL:HG11	2.01	0.42
2:F:49:ALA:HB1	2:F:69:VAL:CG2	2.49	0.42
1:E:195:GLU:HG3	1:E:206:VAL:HG22	2.01	0.42
1:L:21:ILE:HD12	1:L:73:PHE:HD2	1.84	0.42
1:A:160:LEU:HB2	1:A:178:THR:HG22	2.00	0.42
1:A:212:ASN:HD22	1:C:30:SER:HB3	1.84	0.42
2:B:22:CYS:HB3	2:B:78:LEU:HB3	2.01	0.42
2:B:49:ALA:HB1	2:B:69:VAL:CG2	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:VAL:HA	1:A:178:THR:O	2.20	0.42
1:E:146:VAL:HG22	1:E:196:VAL:HG22	2.02	0.42
1:E:23:CYS:HB2	1:E:35:TRP:CH2	2.55	0.42
1:L:35:TRP:HB2	1:L:48:ILE:HB	2.01	0.42
2:B:40:ALA:HB3	2:B:43:ARG:HB3	2.00	0.41
2:D:84:SER:HA	2:D:111:VAL:HB	2.02	0.41
2:F:84:SER:HA	2:F:111:VAL:HB	2.02	0.41
1:A:21:ILE:HD12	1:A:73:PHE:HD2	1.85	0.41
1:E:21:ILE:HD12	1:E:73:PHE:HD2	1.85	0.41
1:A:193:THR:HG23	1:A:206:VAL:HG13	2.03	0.41
2:D:105:GLN:H	2:D:105:GLN:HG3	1.51	0.41
1:E:163:VAL:HG22	1:E:175:LEU:HD13	2.03	0.41
1:C:21:ILE:HD12	1:C:73:PHE:HD2	1.85	0.41
2:D:49:ALA:HB1	2:D:69:VAL:HG21	2.01	0.41
1:A:23:CYS:HB2	1:A:35:TRP:CH2	2.56	0.41
1:L:6:GLN:HG3	1:L:100:PRO:HD2	2.02	0.41
2:D:194:PRO:HD2	2:D:199:THR:HG21	2.03	0.40
1:E:12:PRO:HA	1:E:105:GLU:HG3	2.02	0.40
2:B:105:GLN:H	2:B:105:GLN:HG3	1.51	0.40
1:C:83:ILE:HG13	1:C:106:ILE:HG12	2.04	0.40
1:E:94:TYR:HB3	2:F:58:LYS:HD3	2.03	0.40
2:F:194:PRO:HD2	2:F:199:THR:HG21	2.04	0.40
1:L:32:TYR:O	1:L:90:GLN: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	211/213 (99%)	206 (98%)	5 (2%)	0	100	100
1	C	211/213 (99%)	206 (98%)	5 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	211/213 (99%)	206 (98%)	5 (2%)	0	100	100
1	L	211/213 (99%)	207 (98%)	4 (2%)	0	100	100
2	B	206/216 (95%)	203 (98%)	3 (2%)	0	100	100
2	D	207/216 (96%)	201 (97%)	6 (3%)	0	100	100
2	F	207/216 (96%)	202 (98%)	5 (2%)	0	100	100
2	H	207/216 (96%)	202 (98%)	4 (2%)	1 (0%)	24	43
All	All	1671/1716 (97%)	1633 (98%)	37 (2%)	1 (0%)	48	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	135	THR

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	190/190 (100%)	180 (95%)	10 (5%)	20	42
1	C	190/190 (100%)	183 (96%)	7 (4%)	30	57
1	E	190/190 (100%)	177 (93%)	13 (7%)	14	31
1	L	190/190 (100%)	183 (96%)	7 (4%)	30	57
2	B	178/184 (97%)	164 (92%)	14 (8%)	11	24
2	D	179/184 (97%)	168 (94%)	11 (6%)	17	35
2	F	179/184 (97%)	159 (89%)	20 (11%)	6	12
2	H	179/184 (97%)	157 (88%)	22 (12%)	4	10
All	All	1475/1496 (99%)	1371 (93%)	104 (7%)	13	29

All (104) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	THR

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Mol	Chain	Res	Type
1	A	7	SER
1	A	52	ASN
1	A	89	GLN
1	A	93	ASN
1	A	105	GLU
1	A	153	SER
1	A	178	THR
1	A	181	LEU
1	A	197	THR
2	B	3	GLN
2	B	11	LEU
2	B	12	VAL
2	B	18	LEU
2	B	55	ILE
2	B	56	ASN
2	B	69	VAL
2	B	94	ARG
2	B	97	TRP
2	B	105	GLN
2	B	107	THR
2	B	156	ILE
2	B	218	THR
2	B	226	GLU
1	C	1	ASP
1	C	5	THR
1	C	52	ASN
1	C	93	ASN
1	C	108	ARG
1	C	178	THR
1	C	199	LYS
2	D	3	GLN
2	D	20	LEU
2	D	55	ILE
2	D	68	THR
2	D	69	VAL
2	D	94	ARG
2	D	105	GLN
2	D	107	THR
2	D	199	THR
2	D	205	THR
2	D	226	GLU
1	E	1	ASP

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Mol	Chain	Res	Type
1	E	5	THR
1	E	7	SER
1	E	28	ASP
1	E	52	ASN
1	E	54	LEU
1	E	71	SER
1	E	89	GLN
1	E	93	ASN
1	E	105	GLU
1	E	197	THR
1	E	199	LYS
1	E	212	ASN
2	F	1	GLU
2	F	3	GLN
2	F	11	LEU
2	F	18	LEU
2	F	20	LEU
2	F	21	SER
2	F	56	ASN
2	F	69	VAL
2	F	82	MET
2	F	82(A)	ASN
2	F	82(B)	ILE
2	F	82(C)	LEU
2	F	94	ARG
2	F	97	TRP
2	F	105	GLN
2	F	107	THR
2	F	156	ILE
2	F	199	THR
2	F	218	THR
2	F	226	GLU
2	H	3	GLN
2	H	18	LEU
2	H	20	LEU
2	H	21	SER
2	H	55	ILE
2	H	56	ASN
2	H	68	THR
2	H	69	VAL
2	H	82	MET
2	H	82(A)	ASN

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Mol	Chain	Res	Type
2	H	82(B)	ILE
2	H	82(C)	LEU
2	H	94	ARG
2	H	105	GLN
2	H	107	THR
2	H	116	THR
2	H	135	THR
2	H	180	SER
2	H	186	SER
2	H	199	THR
2	H	205	THR
2	H	218	THR
1	L	7	SER
1	L	24	GLN
1	L	54	LEU
1	L	93	ASN
1	L	112	LYS
1	L	153	SER
1	L	199	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	ASN
1	A	89	GLN
1	A	189	HIS
1	A	190	ASN
2	B	101	ASN
2	B	105	GLN
2	B	212	HIS
1	C	38	GLN
1	C	52	ASN
1	C	137	ASN
1	C	190	ASN
2	D	3	GLN
2	D	39	GLN
2	D	56	ASN
2	D	105	GLN
2	D	204	GLN
1	E	34	ASN
1	E	52	ASN
1	E	89	GLN

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Mol	Chain	Res	Type
1	E	212	ASN
2	F	3	GLN
2	F	39	GLN
2	F	105	GLN
2	F	204	GLN
2	H	39	GLN
2	H	81	GLN
2	H	105	GLN
2	H	204	GLN
1	L	22	ASN
1	L	38	GLN
1	L	137	ASN
1	L	161	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

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	213/213 (100%)	0.06	1 (0%) 87 85	32, 59, 102, 124	0
1	C	213/213 (100%)	-0.03	1 (0%) 87 85	36, 54, 86, 105	0
1	E	213/213 (100%)	0.10	1 (0%) 87 85	39, 57, 91, 117	0
1	L	213/213 (100%)	-0.02	1 (0%) 87 85	35, 58, 87, 100	0
2	B	210/216 (97%)	0.24	4 (1%) 66 62	33, 63, 96, 109	0
2	D	211/216 (97%)	0.10	3 (1%) 73 70	38, 55, 87, 104	0
2	F	211/216 (97%)	0.12	3 (1%) 73 70	41, 56, 87, 105	0
2	H	211/216 (97%)	0.19	6 (2%) 55 50	38, 60, 88, 103	0
All	All	1695/1716 (98%)	0.09	20 (1%) 76 73	32, 58, 92, 124	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	127	ALA	3.4
2	H	127	ALA	3.2
2	H	97	TRP	3.2
2	H	128	CYS	2.9
2	B	134	THR	2.8
2	B	97	TRP	2.6
2	F	142	CYS	2.5
2	H	134	THR	2.5
1	E	23	CYS	2.5
2	F	97	TRP	2.5
2	H	42	GLY	2.4
2	H	142	CYS	2.4
2	D	128	CYS	2.4
1	C	5	THR	2.3
2	D	134	THR	2.3
1	A	94	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
2	F	135	THR	2.3
2	B	55	ILE	2.2
2	D	97	TRP	2.2
1	L	21	ILE	2.0

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.