



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 06:32 AM UTC

PDB ID : 3R08 / pdb_00003r08
Title : Crystal structure of mouse cd3epsilon in complex with antibody 2C11 Fab
Authors : Shore, D.A.; Zhu, X.; Wilson, I.A.
Deposited on : 2011-03-07
Resolution : 4.10 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

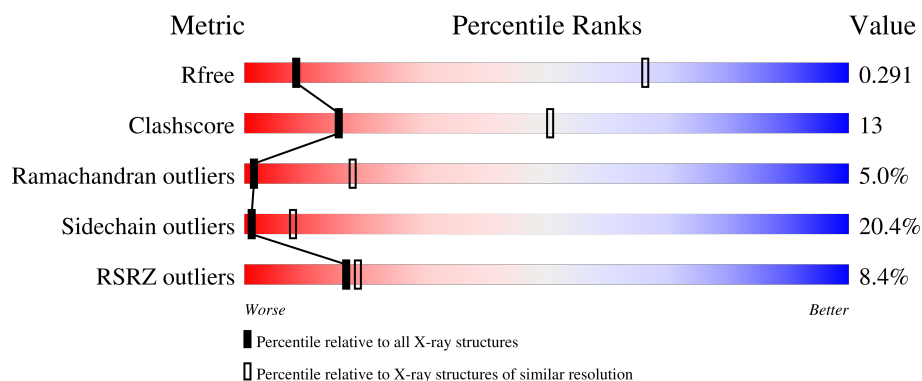
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.10 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	1243 (4.40-3.80)
Clashscore	190562	1293 (4.40-3.80)
Ramachandran outliers	187476	1206 (4.40-3.80)
Sidechain outliers	187428	1193 (4.40-3.80)
RSRZ outliers	180081	1240 (4.40-3.80)

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	L	213	7% 50% 38% 10% .
2	H	216	8% 53% 38% 7% .
3	E	82	13% 50% 39% 7% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3944 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 Mouse anti-mouse CD3epsilon antibody 2C11 light chain.

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	216	Total	C	N	O	S	0	0	0
			1626	1029	268	320	9			

- Molecule 3 is a protein called T-cell surface glycoprotein CD3 epsilon chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	82	Total	C	N	O	S	0	0	0
			654	410	105	137	2			

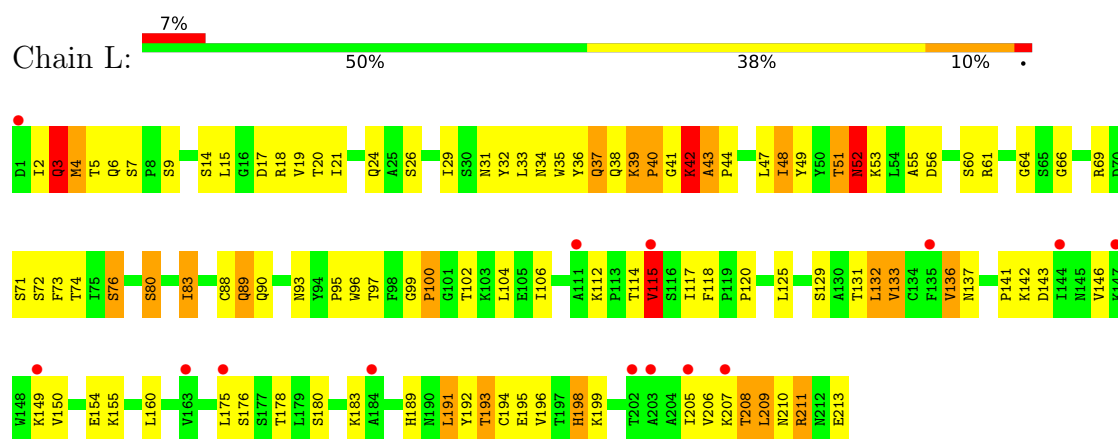
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	80	SER	-	expression tag	UNP P22646
E	81	GLU	-	expression tag	UNP P22646
E	82	TYR	-	expression tag	UNP P22646

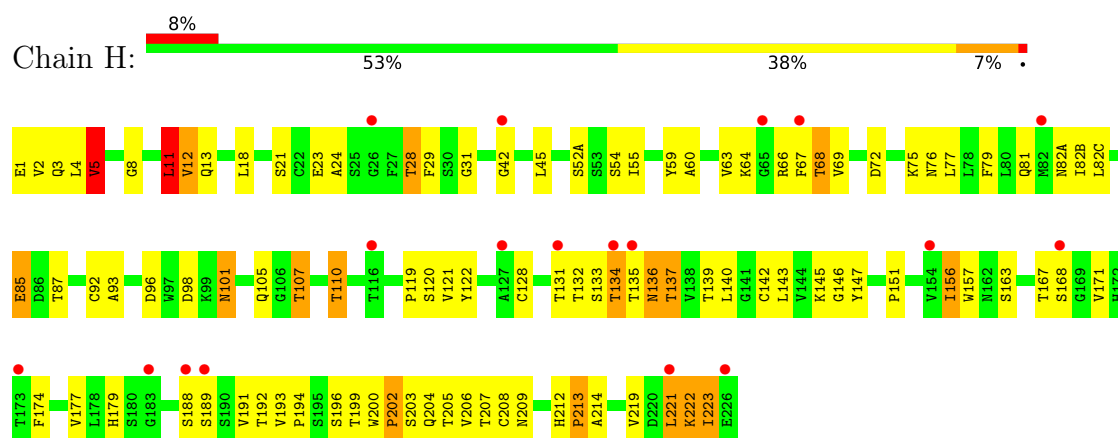
3 Residue-property plots [i](#)

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.

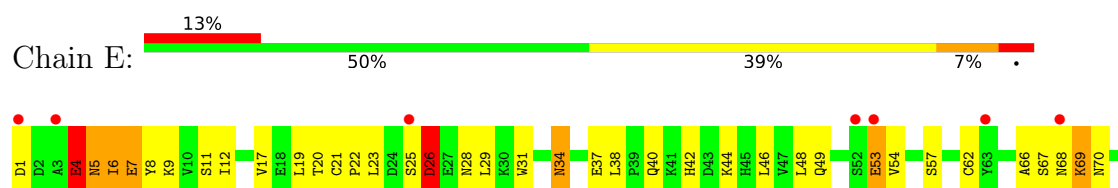
- Molecule 1: Mouse anti-mouse CD3epsilon antibody 2C11 light chain

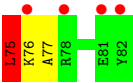


- Molecule 2: Mouse anti-mouse CD3epsilon antibody 2C11 heavy chain



- Molecule 3: T-cell surface glycoprotein CD3 epsilon chain





4 Data and refinement statistics

Property	Value	Source
Space group	I 41 3 2	Depositor
Cell constants a, b, c, α , β , γ	263.18Å 263.18Å 263.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.15 – 4.10 33.15 – 4.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) (33.15-4.10) 96.6 (33.15-4.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 4.13Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, R_{free}	0.222 , 0.277 0.239 , 0.291	Depositor DCC
R_{free} test set	584 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	158.3	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 254.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	3944	wwPDB-VP
Average B, all atoms (Å ²)	180.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.64% 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	L	0.87	1/1701 (0.1%)	1.47	20/2309 (0.9%)
2	H	0.85	0/1666	1.45	22/2274 (1.0%)
3	E	0.95	0/667	1.64	11/904 (1.2%)
All	All	0.88	1/4034 (0.0%)	1.49	53/5487 (1.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	39	LYS	CA-C	5.53	1.59	1.52

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	143	ASP	N-CA-C	8.69	121.66	109.15
1	L	52	ASN	CA-CB-CG	7.83	120.43	112.60
3	E	70	ASN	CA-C-N	7.17	132.67	121.99
3	E	70	ASN	C-N-CA	7.17	132.67	121.99
2	H	5	VAL	N-CA-C	6.83	118.77	106.61
1	L	40	PRO	N-CA-C	6.78	126.43	112.47
1	L	141	PRO	CA-C-N	6.69	129.57	120.54
1	L	141	PRO	C-N-CA	6.69	129.57	120.54
1	L	42	LYS	N-CA-C	6.55	119.40	110.55
2	H	96	ASP	CA-CB-CG	6.50	119.10	112.60
3	E	75	LEU	CA-C-N	6.21	133.40	121.54
3	E	75	LEU	C-N-CA	6.21	133.40	121.54
2	H	85	GLU	N-CA-C	-6.13	104.15	111.69
1	L	74	THR	N-CA-C	6.03	118.22	107.80
2	H	221	LEU	N-CA-C	5.97	118.37	109.59
1	L	41	GLY	CA-C-N	5.95	133.82	121.32
1	L	41	GLY	C-N-CA	5.95	133.82	121.32
2	H	107	THR	CB-CA-C	5.94	120.58	109.71
2	H	135	THR	CA-C-N	5.92	132.84	121.54
2	H	135	THR	C-N-CA	5.92	132.84	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	134	THR	CA-C-N	5.90	129.25	120.17
2	H	134	THR	C-N-CA	5.90	129.25	120.17
3	E	8	TYR	N-CA-C	5.89	118.36	110.35
2	H	93	ALA	N-CA-C	5.74	118.01	108.99
2	H	72	ASP	CA-CB-CG	5.70	118.30	112.60
1	L	118	PHE	N-CA-C	5.69	118.51	109.40
3	E	68	ASN	N-CA-C	-5.63	106.73	113.21
1	L	198	HIS	N-CA-C	5.62	117.94	109.23
1	L	56	ASP	N-CA-C	5.62	117.24	109.15
3	E	26	ASP	N-CA-C	5.58	122.68	110.80
2	H	11	LEU	CA-C-N	5.56	129.42	121.96
2	H	11	LEU	C-N-CA	5.56	129.42	121.96
1	L	115	VAL	N-CA-C	5.54	116.35	107.98
2	H	189	SER	N-CA-C	5.53	117.23	107.61
1	L	4	MET	N-CA-C	5.50	118.79	109.76
2	H	42	GLY	CA-C-N	5.50	129.40	120.60
2	H	42	GLY	C-N-CA	5.50	129.40	120.60
1	L	3	GLN	N-CA-C	5.44	117.24	108.32
2	H	207	THR	CB-CA-C	5.42	118.69	109.80
3	E	42	HIS	N-CA-C	5.37	121.38	114.56
1	L	155	LYS	CA-C-N	5.36	127.73	120.65
1	L	155	LYS	C-N-CA	5.36	127.73	120.65
3	E	69	LYS	CA-C-N	5.29	130.33	122.82
3	E	69	LYS	C-N-CA	5.29	130.33	122.82
2	H	21	SER	N-CA-C	5.25	117.58	107.44
1	L	199	LYS	CA-C-N	5.24	127.55	120.38
1	L	199	LYS	C-N-CA	5.24	127.55	120.38
3	E	7	GLU	CB-CG-CD	5.23	121.49	112.60
2	H	156	ILE	N-CA-CB	5.13	117.93	111.46
2	H	202	PRO	CA-C-N	5.11	128.48	120.82
2	H	202	PRO	C-N-CA	5.11	128.48	120.82
2	H	68	THR	N-CA-C	5.08	115.72	107.23
1	L	133	VAL	N-CA-C	5.04	116.98	108.81

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	L	1664	0	1608	52	0
2	H	1626	0	1593	42	0
3	E	654	0	612	14	0
All	All	3944	0	3813	104	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 13.

All (104) 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:H:156:ILE:HG13	2:H:209:ASN:HB2	1.62	0.82
2:H:151:PRO:HD2	2:H:213:PRO:HG3	1.66	0.77
2:H:119:PRO:HB3	2:H:147:TYR:HB3	1.67	0.75
2:H:133:SER:HB2	2:H:136:ASN:HA	1.70	0.74
3:E:9:LYS:HB3	3:E:20:THR:HB	1.72	0.72
1:L:34:ASN:HD22	1:L:49:TYR:HA	1.56	0.71
2:H:171:VAL:HG22	2:H:191:VAL:HG23	1.71	0.71
1:L:115:VAL:HG21	1:L:196:VAL:HG21	1.74	0.68
2:H:206:VAL:H	2:H:222:LYS:HE3	1.59	0.68
1:L:90:GLN:HE21	1:L:97:THR:H	1.47	0.62
1:L:189:HIS:O	1:L:211:ARG:NH1	2.32	0.62
1:L:4:MET:HB2	1:L:99:GLY:HA2	1.82	0.62
1:L:120:PRO:HD3	1:L:132:LEU:HG	1.80	0.61
1:L:193:THR:HG23	1:L:208:THR:HG23	1.81	0.61
1:L:52:ASN:H	1:L:52:ASN:HD22	1.48	0.60
2:H:68:THR:HG23	2:H:81:GLN:HB3	1.84	0.60
2:H:5:VAL:HG23	2:H:23:GLU:HB2	1.84	0.60
2:H:87:THR:HG23	2:H:110:THR:HA	1.84	0.59
1:L:83:ILE:HG12	1:L:106:ILE:HB	1.86	0.58
1:L:66:GLY:HA3	1:L:71:SER:HA	1.85	0.58
1:L:15:LEU:HD11	1:L:106:ILE:HD11	1.86	0.57
1:L:36:TYR:OH	1:L:89:GLN:NE2	2.38	0.57
1:L:195:GLU:HA	1:L:206:VAL:HG12	1.86	0.56
1:L:115:VAL:CG2	1:L:196:VAL:HG21	2.34	0.56
3:E:7:GLU:HA	3:E:22:PRO:HB3	1.86	0.56
2:H:66:ARG:HD3	2:H:82(B):ILE:O	2.04	0.56
1:L:2:ILE:HG22	1:L:3:GLN:N	2.20	0.56
1:L:21:ILE:HG12	1:L:102:THR:HG21	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:95:PRO:HG3	2:H:60:ALA:HA	1.89	0.55
2:H:12:VAL:HG11	2:H:82(C):LEU:HD13	1.89	0.54
2:H:212:HIS:HE1	2:H:214:ALA:HB3	1.71	0.54
2:H:29:PHE:HB3	2:H:76:ASN:HD22	1.72	0.54
1:L:37:GLN:HB2	1:L:47:LEU:HD21	1.89	0.53
3:E:29:LEU:HD11	3:E:62:CYS:SG	2.47	0.53
1:L:6:GLN:HB2	1:L:100:PRO:HD2	1.88	0.53
1:L:34:ASN:ND2	1:L:49:TYR:HA	2.23	0.53
3:E:75:LEU:O	3:E:76:LYS:HG2	2.09	0.53
1:L:137:ASN:HD21	2:H:174:PHE:HZ	1.57	0.53
1:L:42:LYS:HE3	1:L:42:LYS:H	1.74	0.52
1:L:136:VAL:HG23	1:L:175:LEU:HB3	1.90	0.52
1:L:146:VAL:HG12	1:L:196:VAL:HA	1.91	0.52
2:H:133:SER:CB	2:H:136:ASN:HA	2.40	0.52
2:H:66:ARG:HG3	2:H:67:PHE:CD1	2.46	0.52
2:H:31:GLY:HA2	2:H:52(A):SER:HB2	1.92	0.51
2:H:66:ARG:HG3	2:H:67:PHE:HD1	1.74	0.51
2:H:212:HIS:CE1	2:H:214:ALA:HB3	2.46	0.51
3:E:67:SER:C	3:E:69:LYS:H	2.19	0.51
1:L:6:GLN:NE2	1:L:102:THR:H	2.09	0.50
1:L:19:VAL:HG11	1:L:104:LEU:HD11	1.93	0.50
2:H:11:LEU:HG	2:H:12:VAL:H	1.77	0.50
1:L:2:ILE:HG22	1:L:3:GLN:H	1.75	0.50
2:H:75:LYS:HB2	2:H:77:LEU:HD12	1.94	0.48
3:E:53:GLU:HG2	3:E:76:LYS:HB2	1.95	0.48
1:L:35:TRP:CZ3	1:L:88:CYS:HB3	2.49	0.48
1:L:5:THR:HB	1:L:24:GLN:HB3	1.95	0.47
1:L:37:GLN:HB2	1:L:47:LEU:HD11	1.97	0.47
1:L:150:VAL:HG22	1:L:192:TYR:HD1	1.80	0.46
2:H:157:TRP:CH2	2:H:208:CYS:HB2	2.51	0.46
2:H:120:SER:HB3	2:H:122:TYR:CZ	2.51	0.45
2:H:82(A):ASN:HB2	2:H:82(B):ILE:HD12	1.98	0.45
2:H:142:CYS:HB2	2:H:157:TRP:CZ2	2.52	0.45
1:L:43:ALA:HA	1:L:44:PRO:HD3	1.88	0.45
1:L:48:ILE:HG22	1:L:53:LYS:O	2.16	0.45
3:E:21:CYS:HB2	3:E:31:TRP:CZ2	2.52	0.45
1:L:48:ILE:HD12	1:L:64:GLY:N	2.32	0.45
1:L:131:THR:HG23	1:L:180:SER:HA	1.99	0.44
1:L:150:VAL:HG22	1:L:192:TYR:CD1	2.52	0.44
3:E:21:CYS:SG	3:E:22:PRO:HD2	2.58	0.44
1:L:49:TYR:HE1	1:L:55:ALA:HA	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:4:LEU:HD23	2:H:24:ALA:HA	1.99	0.44
2:H:59:TYR:HE1	2:H:69:VAL:HG12	1.81	0.44
3:E:12:ILE:HG13	3:E:17:VAL:HG13	2.00	0.44
1:L:209:LEU:HG	2:H:128:CYS:HB2	2.00	0.44
2:H:120:SER:HB2	2:H:145:LYS:HB3	1.99	0.44
2:H:121:VAL:HB	2:H:219:VAL:HG11	1.99	0.43
1:L:150:VAL:HG13	1:L:192:TYR:CE1	2.53	0.43
1:L:196:VAL:HB	1:L:205:ILE:O	2.19	0.43
2:H:200:TRP:CD1	2:H:202:PRO:HA	2.54	0.43
2:H:140:LEU:HD22	2:H:223:ILE:HD13	2.01	0.43
3:E:46:LEU:HD23	3:E:48:LEU:HD11	1.99	0.43
2:H:156:ILE:HD12	2:H:163:SER:HA	2.01	0.42
1:L:117:ILE:HD13	1:L:194:CYS:HB2	2.02	0.42
1:L:29:ILE:HD12	1:L:31:ASN:H	1.84	0.42
2:H:63:VAL:O	2:H:64:LYS:C	2.63	0.42
3:E:57:SER:OG	3:E:76:LYS:HB3	2.20	0.42
2:H:133:SER:HB2	2:H:136:ASN:CA	2.44	0.41
2:H:200:TRP:CG	2:H:202:PRO:HA	2.55	0.41
1:L:61:ARG:HB2	1:L:76:SER:HB3	2.02	0.41
3:E:4:GLU:HB3	3:E:5:ASN:H	1.69	0.41
1:L:2:ILE:CG2	1:L:3:GLN:N	2.84	0.41
1:L:52:ASN:H	1:L:52:ASN:ND2	2.15	0.41
1:L:125:LEU:O	1:L:183:LYS:HD2	2.21	0.41
1:L:160:LEU:HD21	2:H:179:HIS:HB2	2.01	0.41
2:H:157:TRP:CZ3	2:H:208:CYS:HB2	2.55	0.41
1:L:191:LEU:HD12	1:L:210:ASN:HB2	2.02	0.41
2:H:137:THR:HA	2:H:194:PRO:HA	2.03	0.41
3:E:26:ASP:C	3:E:28:ASN:H	2.29	0.41
3:E:31:TRP:HB3	3:E:38:LEU:HD12	2.03	0.40
1:L:193:THR:HG22	1:L:206:VAL:HB	2.04	0.40
1:L:48:ILE:HD11	1:L:73:PHE:HE1	1.87	0.40
2:H:199:THR:HA	2:H:204:GLN:NE2	2.37	0.40
1:L:2:ILE:CG2	1:L:3:GLN:H	2.34	0.40
1:L:207:LYS:HD2	1:L:207:LYS:HA	2.01	0.40
2:H:101:ASN:HD22	2:H:101:ASN:HA	1.52	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	L	211/213 (99%)	179 (85%)	22 (10%)	10 (5%)	2	19
2	H	214/216 (99%)	180 (84%)	27 (13%)	7 (3%)	3	24
3	E	80/82 (98%)	61 (76%)	11 (14%)	8 (10%)	0	8
All	All	505/511 (99%)	420 (83%)	60 (12%)	25 (5%)	1	18

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	18	ARG
1	L	40	PRO
1	L	43	ALA
1	L	51	THR
2	H	132	THR
2	H	136	ASN
3	E	26	ASP
1	L	60	SER
1	L	80	SER
1	L	211	ARG
2	H	146	GLY
1	L	26	SER
1	L	32	TYR
1	L	100	PRO
2	H	28	THR
3	E	6	ILE
3	E	77	ALA
2	H	131	THR
3	E	34	ASN
3	E	66	ALA
2	H	213	PRO
3	E	4	GLU
3	E	23	LEU
3	E	40	GLN

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Mol	Chain	Res	Type
2	H	8	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	190/190 (100%)	150 (79%)	40 (21%)	1	7
2	H	184/184 (100%)	148 (80%)	36 (20%)	1	9
3	E	73/76 (96%)	58 (80%)	15 (20%)	1	8
All	All	447/450 (99%)	356 (80%)	91 (20%)	1	8

All (91) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	3	GLN
1	L	7	SER
1	L	9	SER
1	L	14	SER
1	L	17	ASP
1	L	20	THR
1	L	33	LEU
1	L	37	GLN
1	L	38	GLN
1	L	39	LYS
1	L	42	LYS
1	L	48	ILE
1	L	51	THR
1	L	52	ASN
1	L	69	ARG
1	L	72	SER
1	L	76	SER
1	L	80	SER
1	L	83	ILE
1	L	89	GLN
1	L	93	ASN

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Mol	Chain	Res	Type
1	L	96	TRP
1	L	112	LYS
1	L	114	THR
1	L	115	VAL
1	L	129	SER
1	L	132	LEU
1	L	133	VAL
1	L	136	VAL
1	L	142	LYS
1	L	149	LYS
1	L	154	GLU
1	L	176	SER
1	L	178	THR
1	L	191	LEU
1	L	193	THR
1	L	198	HIS
1	L	208	THR
1	L	209	LEU
1	L	213	GLU
2	H	1	GLU
2	H	2	VAL
2	H	3	GLN
2	H	5	VAL
2	H	11	LEU
2	H	12	VAL
2	H	13	GLN
2	H	18	LEU
2	H	28	THR
2	H	45	LEU
2	H	54	SER
2	H	55	ILE
2	H	79	PHE
2	H	85	GLU
2	H	92	CYS
2	H	98	ASP
2	H	101	ASN
2	H	105	GLN
2	H	107	THR
2	H	110	THR
2	H	134	THR
2	H	137	THR
2	H	139	THR

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Mol	Chain	Res	Type
2	H	143	LEU
2	H	167	THR
2	H	168	SER
2	H	177	VAL
2	H	188	SER
2	H	192	THR
2	H	193	VAL
2	H	196	SER
2	H	203	SER
2	H	205	THR
2	H	221	LEU
2	H	222	LYS
2	H	223	ILE
3	E	1	ASP
3	E	4	GLU
3	E	5	ASN
3	E	6	ILE
3	E	11	SER
3	E	19	LEU
3	E	25	SER
3	E	26	ASP
3	E	34	ASN
3	E	37	GLU
3	E	44	LYS
3	E	49	GLN
3	E	53	GLU
3	E	54	VAL
3	E	75	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	38	GLN
1	L	52	ASN
1	L	89	GLN
1	L	138	ASN
1	L	145	ASN
1	L	161	GLN
1	L	190	ASN
1	L	212	ASN
2	H	3	GLN
2	H	39	GLN

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Mol	Chain	Res	Type
2	H	76	ASN
2	H	81	GLN
2	H	101	ASN
2	H	136	ASN
2	H	179	HIS
2	H	204	GLN
3	E	28	ASN

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	L	213/213 (100%)	0.79	14 (6%)	24 23	131, 169, 223, 232	0
2	H	216/216 (100%)	0.71	18 (8%)	17 19	151, 182, 213, 238	0
3	E	82/82 (100%)	0.94	11 (13%)	7 11	166, 180, 190, 215	0
All	All	511/511 (100%)	0.78	43 (8%)	17 19	131, 180, 221, 238	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	E	1	ASP	5.6
2	H	134	THR	5.1
1	L	202	THR	4.1
2	H	188	SER	3.6
2	H	65	GLY	3.5
1	L	144	ILE	3.2
1	L	207	LYS	3.2
2	H	226	GLU	3.1
1	L	184	ALA	3.1
2	H	82	MET	3.0
3	E	81	GLU	2.9
2	H	173	THR	2.9
2	H	221	LEU	2.8
3	E	25	SER	2.8
2	H	26	GLY	2.7
2	H	135	THR	2.7
3	E	78	ARG	2.6
2	H	131	THR	2.6
3	E	52	SER	2.6
3	E	53	GLU	2.6
2	H	116	THR	2.6
2	H	168	SER	2.5
1	L	1	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	L	205	ILE	2.4
2	H	154	VAL	2.4
2	H	67	PHE	2.4
3	E	63	TYR	2.4
2	H	42	GLY	2.3
3	E	82	TYR	2.3
1	L	147	LYS	2.3
1	L	115	VAL	2.3
3	E	3	ALA	2.2
1	L	175	LEU	2.2
2	H	183	GLY	2.2
2	H	127	ALA	2.1
3	E	68	ASN	2.1
1	L	163	VAL	2.1
1	L	111	ALA	2.1
1	L	135	PHE	2.1
1	L	149	LYS	2.0
3	E	76	LYS	2.0
2	H	189	SER	2.0
1	L	203	ALA	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.