



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 10:28 AM UTC

PDB ID : 2IAL / pdb_00002ial
Title : Structural basis for recognition of mutant self by a tumor-specific, MHC class II-restricted TCR
Authors : Deng, L.; Langley, R.J.; Mariuzza, R.A.
Deposited on : 2006-09-08
Resolution : 1.92 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

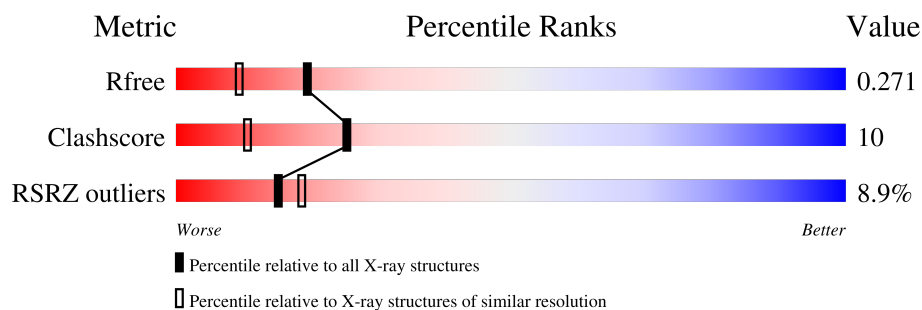
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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	1188 (1.92-1.92)
Clashscore	190562	1209 (1.92-1.92)
RSRZ outliers	180081	1188 (1.92-1.92)

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	202	21% 79% 17% ..
1	C	202	8% 82% 13% ..
2	B	240	3% 79% 19% .
2	D	240	5% 80% 19% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7105 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 CD4+ T cell receptor E8 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	S	0	0	0
			1534	961	258	308	7			
1	C	195	Total	C	N	O	S	0	0	0
			1526	957	257	305	7			

- Molecule 2 is a protein called CD4+ T cell receptor E8 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	239	Total	C	N	O	S	0	0	0
			1901	1206	324	362	9			
2	D	239	Total	C	N	O	S	0	0	0
			1901	1206	324	362	9			

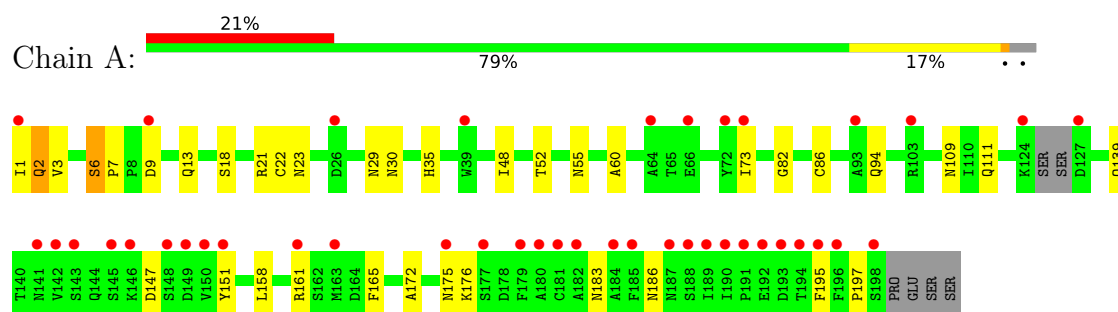
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	27	Total	O	0	0
			27	27		
3	B	94	Total	O	0	0
			94	94		
3	C	44	Total	O	0	0
			44	44		
3	D	78	Total	O	0	0
			78	78		

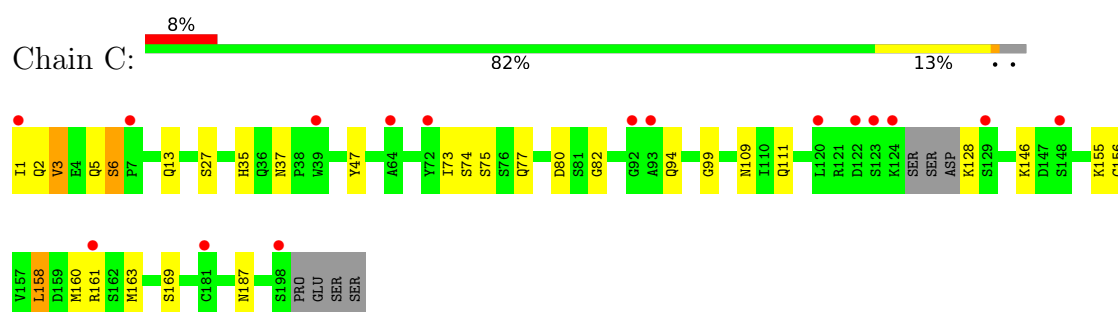
3 Residue-property plots [i](#)

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.

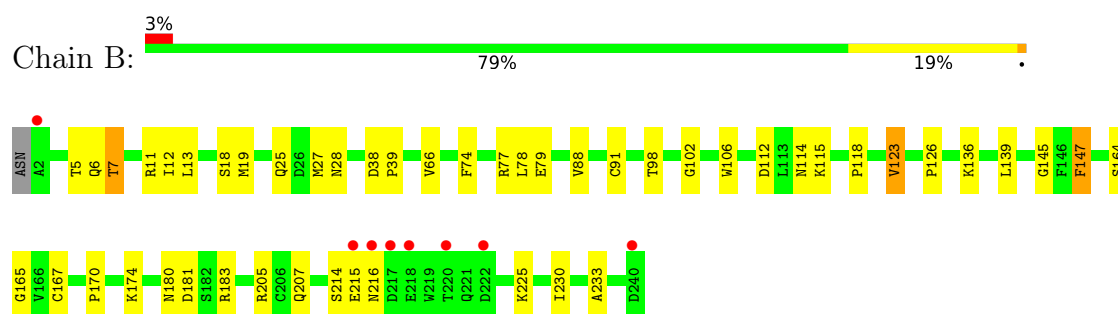
- Molecule 1: CD4+ T cell receptor E8 alpha chain



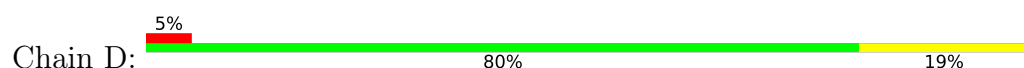
- Molecule 1: CD4+ T cell receptor E8 alpha chain

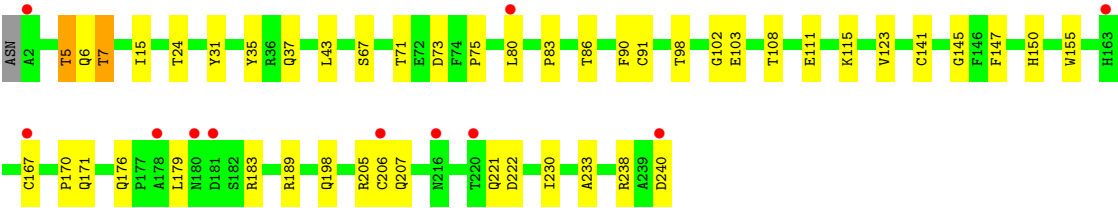


- Molecule 2: CD4+ T cell receptor E8 beta chain



- Molecule 2: CD4+ T cell receptor E8 beta chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	170.04Å 65.66Å 84.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.31 – 1.92 44.31 – 1.92	Depositor EDS
% Data completeness (in resolution range)	94.1 (44.31-1.92) 94.3 (44.31-1.92)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.03 (at 1.92Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.208 , 0.265 (Not available) , 0.271	Depositor DCC
R_{free} test set	3495 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.178	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 29.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7105	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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	0.95	1/1566 (0.1%)	1.08	5/2129 (0.2%)
1	C	1.07	1/1558 (0.1%)	1.22	8/2118 (0.4%)
2	B	1.18	2/1956 (0.1%)	1.19	6/2666 (0.2%)
2	D	1.08	2/1956 (0.1%)	1.20	2/2666 (0.1%)
All	All	1.08	6/7036 (0.1%)	1.17	21/9579 (0.2%)

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	2
1	C	0	1
2	B	0	4
2	D	0	3
All	All	0	10

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	123	VAL	CA-CB	6.05	1.62	1.54
2	D	5	THR	CA-CB	6.04	1.61	1.53
1	C	73	ILE	CA-CB	6.03	1.61	1.54
2	B	118	PRO	CA-C	5.62	1.57	1.52
1	A	48	ILE	CA-CB	5.40	1.61	1.54
2	D	71	THR	CA-CB	5.20	1.61	1.53

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	7	THR	CA-C-N	16.62	140.61	119.84
2	D	7	THR	C-N-CA	16.62	140.61	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	147	PHE	CA-C-N	15.73	139.50	119.84
2	B	147	PHE	C-N-CA	15.73	139.50	119.84
1	C	6	SER	CA-C-N	15.66	136.51	120.38
1	C	6	SER	C-N-CA	15.66	136.51	120.38
1	A	7	PRO	N-CA-C	-9.12	99.58	110.70
2	B	7	THR	CA-C-N	7.49	129.21	119.84
2	B	7	THR	C-N-CA	7.49	129.21	119.84
1	C	158	LEU	CA-CB-CG	6.02	137.37	116.30
1	A	6	SER	CA-C-N	5.95	126.51	120.38
1	A	6	SER	C-N-CA	5.95	126.51	120.38
1	C	2	GLN	N-CA-C	5.87	117.70	108.96
1	C	156	CYS	CA-CB-SG	5.82	127.79	114.40
2	B	25	GLN	N-CA-CB	-5.62	101.61	110.77
1	A	2	GLN	N-CA-C	5.48	117.72	109.23
1	A	94	GLN	N-CA-C	-5.47	104.81	112.25
1	C	3	VAL	N-CA-C	-5.10	101.29	108.27
2	B	216	ASN	N-CA-C	-5.06	107.16	113.38
1	C	37	ASN	CA-C-N	5.01	126.10	119.84
1	C	37	ASN	C-N-CA	5.01	126.10	119.84

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	6	SER	Peptide,Mainchain
2	B	147	PHE	Peptide,Mainchain
2	B	7	THR	Peptide,Mainchain
1	C	6	SER	Peptide
2	D	147	PHE	Peptide,Mainchain
2	D	7	THR	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	1534	0	1463	38	0
1	C	1526	0	1459	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1901	0	1800	41	0
2	D	1901	0	1800	40	0
3	A	27	0	0	0	0
3	B	94	0	0	6	0
3	C	44	0	0	2	0
3	D	78	0	0	6	0
All	All	7105	0	6522	128	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 10.

All (128) 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:175:ASN:ND2	1:A:175:ASN:H	1.75	0.84
1:A:21:ARG:HE	1:A:23:ASN:HD21	1.22	0.83
1:A:55:ASN:HD21	1:C:161:ARG:HH22	1.24	0.83
2:B:77:ARG:HD3	2:B:79:GLU:OE2	1.79	0.83
2:B:170:PRO:HD2	3:B:322:HOH:O	1.80	0.81
2:B:180:ASN:HB3	3:B:307:HOH:O	1.79	0.81
2:B:112:ASP:HB2	2:B:114:ASN:ND2	1.95	0.81
1:C:111:GLN:CD	1:C:111:GLN:H	1.90	0.80
2:B:98:THR:CG2	3:B:321:HOH:O	2.31	0.78
2:B:5:THR:HG23	3:B:245:HOH:O	1.86	0.76
1:C:160:MET:HE2	1:C:163:MET:HE3	1.69	0.73
1:A:55:ASN:ND2	1:C:161:ARG:HH22	1.86	0.73
1:A:147:ASP:OD2	1:A:176:LYS:NZ	2.22	0.72
2:D:123:VAL:HG23	2:D:233:ALA:HB3	1.72	0.72
1:A:21:ARG:HE	1:A:23:ASN:ND2	1.89	0.69
1:A:183:ASN:HA	1:A:186:ASN:OD1	1.92	0.69
2:B:112:ASP:HB2	2:B:114:ASN:HD21	1.57	0.69
2:D:238:ARG:HG2	2:D:240:ASP:HB2	1.75	0.68
2:B:115:LYS:HD3	2:B:181:ASP:OD1	1.94	0.68
2:B:215:GLU:H	2:B:215:GLU:CD	2.03	0.66
1:A:13:GLN:HB3	1:A:109:ASN:HD21	1.60	0.66
1:C:111:GLN:H	1:C:111:GLN:NE2	1.94	0.65
2:D:24:THR:HG23	2:D:73:ASP:OD1	1.97	0.64
2:B:114:ASN:H	2:B:114:ASN:HD22	1.44	0.62
2:D:221:GLN:HG3	2:D:222:ASP:N	2.15	0.62
1:C:3:VAL:HB	3:C:244:HOH:O	1.99	0.62
2:D:170:PRO:HD2	3:D:301:HOH:O	1.98	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:5:GLN:HE21	1:C:99:GLY:HA3	1.66	0.61
2:D:171:GLN:H	2:D:171:GLN:HE21	1.47	0.60
1:A:175:ASN:H	1:A:175:ASN:HD22	1.48	0.60
2:D:150:HIS:HD2	3:D:298:HOH:O	1.83	0.59
1:A:3:VAL:CG1	1:A:86:CYS:SG	2.90	0.59
1:A:161:ARG:HE	2:D:205:ARG:HH21	1.50	0.58
1:A:161:ARG:HE	2:D:205:ARG:NH2	2.01	0.58
2:D:221:GLN:HG3	2:D:222:ASP:H	1.68	0.57
1:C:13:GLN:HE21	1:C:109:ASN:ND2	2.03	0.57
2:D:6:GLN:NE2	2:D:91:CYS:H	2.02	0.57
2:B:6:GLN:HE22	2:B:91:CYS:H	1.51	0.56
2:D:145:GLY:O	2:D:183:ARG:HD3	2.04	0.56
1:C:158:LEU:HD22	3:C:242:HOH:O	2.04	0.56
2:B:98:THR:HG23	3:B:321:HOH:O	1.97	0.56
2:B:115:LYS:CD	2:B:181:ASP:OD1	2.54	0.56
2:D:5:THR:HG23	3:D:303:HOH:O	2.05	0.56
1:A:175:ASN:ND2	1:A:175:ASN:N	2.51	0.55
1:C:169:SER:OG	2:D:189:ARG:HD3	2.06	0.55
1:A:13:GLN:HE21	1:A:109:ASN:HD22	1.55	0.55
2:B:114:ASN:ND2	2:B:114:ASN:H	2.03	0.55
2:D:6:GLN:HE21	2:D:102:GLY:HA3	1.72	0.54
1:A:158:LEU:HD11	2:B:165:GLY:O	2.07	0.54
1:C:160:MET:CE	1:C:163:MET:HE3	2.38	0.53
2:D:6:GLN:HE22	2:D:90:PHE:HA	1.73	0.53
1:A:158:LEU:HB2	2:B:167:CYS:HB2	1.91	0.53
2:B:145:GLY:O	2:B:183:ARG:HD3	2.09	0.52
2:D:15:ILE:HD11	2:D:83:PRO:HG3	1.91	0.52
2:D:171:GLN:NE2	3:D:301:HOH:O	2.42	0.52
2:B:214:SER:O	2:B:225:LYS:NZ	2.29	0.51
2:D:171:GLN:HE21	2:D:171:GLN:N	2.08	0.51
1:A:195:PHE:CE2	1:A:197:PRO:HG3	2.46	0.51
2:D:67:SER:O	2:D:75:PRO:HD2	2.10	0.51
2:B:66:VAL:HG13	2:B:74:PHE:HE1	1.76	0.50
2:B:6:GLN:HE21	2:B:102:GLY:HA3	1.76	0.50
2:B:207:GLN:HG3	2:B:230:ILE:HG23	1.94	0.50
2:B:6:GLN:NE2	2:B:91:CYS:H	2.10	0.50
1:C:1:ILE:N	1:C:27:SER:OG	2.43	0.50
2:B:174:LYS:HE2	3:B:328:HOH:O	2.12	0.49
1:A:161:ARG:HH21	2:D:205:ARG:HH21	1.59	0.49
1:A:165:PHE:CE2	2:B:136:LYS:HE2	2.47	0.48
2:B:112:ASP:OD2	2:B:114:ASN:ND2	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:ILE:HG13	1:A:2:GLN:N	2.29	0.48
1:C:47:TYR:HB2	2:D:98:THR:HG22	1.95	0.47
1:C:94:GLN:HG3	2:D:31:TYR:CE1	2.49	0.47
1:A:165:PHE:CD2	2:B:136:LYS:HE2	2.50	0.47
1:A:111:GLN:HA	1:A:111:GLN:OE1	2.14	0.47
1:A:139:GLN:H	1:A:139:GLN:HG2	1.56	0.46
2:B:38:ASP:HB3	2:B:39:PRO:HD2	1.97	0.46
2:D:198:GLN:HA	2:D:238:ARG:O	2.16	0.46
1:A:158:LEU:HD23	1:A:158:LEU:C	2.41	0.46
2:B:18:SER:HB2	2:B:77:ARG:NH1	2.31	0.46
2:D:123:VAL:HG23	2:D:233:ALA:CB	2.44	0.45
2:B:145:GLY:O	2:B:183:ARG:CD	2.64	0.45
2:D:176:GLN:OE1	2:D:179:LEU:HD13	2.17	0.45
1:A:29:ASN:OD1	1:A:30:ASN:ND2	2.48	0.45
2:D:238:ARG:HD2	3:D:285:HOH:O	2.17	0.45
1:A:175:ASN:HD22	1:A:175:ASN:N	2.13	0.44
2:D:37:GLN:HB2	2:D:43:LEU:HD23	1.99	0.44
2:B:126:PRO:HD3	2:B:139:LEU:HG	1.98	0.44
2:B:13:LEU:HD11	2:B:19:MET:HB2	2.00	0.44
1:C:35:HIS:CE1	1:C:82:GLY:HA3	2.53	0.44
1:C:74:SER:O	1:C:75:SER:C	2.59	0.44
2:D:86:THR:HG23	2:D:108:THR:HA	2.00	0.44
1:C:111:GLN:CD	1:C:111:GLN:N	2.61	0.44
1:C:158:LEU:HB2	2:D:167:CYS:HB2	2.00	0.44
2:D:35:TYR:HB3	2:D:43:LEU:HD22	2.00	0.43
2:D:207:GLN:HG3	2:D:230:ILE:HG23	2.00	0.43
2:B:77:ARG:CD	2:B:79:GLU:OE2	2.57	0.43
1:C:155:LYS:HA	1:C:169:SER:O	2.18	0.43
1:A:161:ARG:NH1	2:B:164:SER:HB2	2.33	0.43
1:C:146:LYS:HD2	1:C:187:ASN:HD21	1.83	0.43
1:A:13:GLN:HB3	1:A:109:ASN:ND2	2.31	0.43
2:D:80:LEU:N	2:D:80:LEU:HD23	2.34	0.43
2:D:103:GLU:HG2	3:D:302:HOH:O	2.18	0.43
1:A:151:TYR:O	1:A:172:ALA:HA	2.19	0.42
2:D:6:GLN:HE22	2:D:91:CYS:H	1.66	0.42
1:A:52:THR:HA	1:A:60:ALA:O	2.20	0.42
2:D:111:GLU:HG3	2:D:115:LYS:NZ	2.34	0.42
2:B:88:VAL:HG22	2:B:106:TRP:CD1	2.55	0.42
2:B:11:ARG:NH1	2:B:12:ILE:O	2.53	0.42
1:C:35:HIS:HE1	1:C:82:GLY:HA3	1.84	0.42
1:A:183:ASN:OD1	1:A:183:ASN:C	2.63	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:66:VAL:CG1	2:B:74:PHE:HE1	2.33	0.42
1:C:13:GLN:HB3	1:C:109:ASN:HD21	1.85	0.42
2:D:15:ILE:HD11	2:D:83:PRO:HD3	2.02	0.41
1:A:3:VAL:HG13	1:A:22:CYS:SG	2.60	0.41
2:B:205:ARG:NH1	2:B:207:GLN:HB2	2.35	0.41
1:A:1:ILE:HG13	1:A:2:GLN:H	1.86	0.41
2:D:123:VAL:HG22	2:D:206:CYS:SG	2.60	0.41
2:B:27:MET:O	2:B:28:ASN:HB2	2.20	0.41
2:B:123:VAL:HG23	2:B:233:ALA:HB3	2.03	0.41
1:C:77:GLN:O	1:C:80:ASP:HB2	2.21	0.41
1:C:128:LYS:HB2	1:C:128:LYS:HE2	1.84	0.41
1:A:18:SER:HB2	1:A:73:ILE:HB	2.03	0.41
1:A:161:ARG:NE	2:D:205:ARG:HH21	2.17	0.41
1:A:9:ASP:OD1	1:A:9:ASP:C	2.64	0.40
2:B:77:ARG:HG3	2:B:78:LEU:N	2.35	0.40
2:D:141:CYS:HB2	2:D:155:TRP:CZ2	2.55	0.40
1:A:35:HIS:HE1	1:A:82:GLY:HA3	1.87	0.40
1:A:3:VAL:HG11	1:A:86:CYS:SG	2.61	0.40
2:B:215:GLU:CD	2:B:215:GLU:N	2.76	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

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

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	196/202 (97%)	1.20	42 (21%) 2 3	21, 37, 68, 77	0
1	C	195/202 (96%)	0.48	16 (8%) 17 21	17, 28, 43, 57	0
2	B	239/240 (99%)	0.21	8 (3%) 49 54	15, 25, 39, 58	0
2	D	239/240 (99%)	0.45	11 (4%) 37 42	17, 28, 46, 55	0
All	All	869/884 (98%)	0.56	77 (8%) 15 19	15, 29, 51, 77	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	189	ILE	5.9
1	A	1	ILE	4.8
1	A	190	ILE	4.8
1	C	1	ILE	4.7
1	A	187	ASN	4.7
2	D	240	ASP	4.4
1	C	123	SER	4.3
1	A	191	PRO	4.1
1	A	93	ALA	3.7
1	A	185	PHE	3.7
2	D	220	THR	3.6
1	A	193	ASP	3.5
1	C	39	TRP	3.4
1	A	181	CYS	3.4
1	A	39	TRP	3.3
2	D	178	ALA	3.3
1	A	184	ALA	3.2
1	A	142	VAL	3.1
2	B	216	ASN	3.1
1	C	181	CYS	3.1
2	D	2	ALA	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	161	ARG	3.0
1	A	175	ASN	3.0
1	A	124	LYS	2.9
2	B	222	ASP	2.9
1	C	93	ALA	2.9
1	A	149	ASP	2.9
2	B	240	ASP	2.8
1	A	188	SER	2.8
1	A	195	PHE	2.8
1	A	163	MET	2.8
1	C	124	LYS	2.7
2	B	217	ASP	2.7
1	A	182	ALA	2.7
1	A	127	ASP	2.7
2	B	215	GLU	2.7
1	A	196	PHE	2.7
1	A	143	SER	2.6
2	D	163	HIS	2.6
1	A	26	ASP	2.6
1	A	198	SER	2.6
1	C	122	ASP	2.6
1	A	141	ASN	2.6
2	D	181	ASP	2.5
1	A	146	LYS	2.5
1	A	180	ALA	2.4
2	B	2	ALA	2.4
1	A	148	SER	2.4
1	C	129	SER	2.4
1	A	151	TYR	2.4
1	C	72	TYR	2.3
2	D	216	ASN	2.3
1	A	179	PHE	2.3
2	D	206	CYS	2.3
2	B	220	THR	2.3
1	A	64	ALA	2.3
1	C	64	ALA	2.3
1	A	73	ILE	2.3
1	A	150	VAL	2.3
2	D	80	LEU	2.3
1	A	72	TYR	2.2
1	A	177	SER	2.2
1	A	145	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	148	SER	2.2
1	A	9	ASP	2.2
1	C	198	SER	2.2
1	A	194	THR	2.1
1	A	103	ARG	2.1
1	A	192	GLU	2.1
1	C	161	ARG	2.1
1	A	66	GLU	2.1
2	B	218	GLU	2.1
1	C	120	LEU	2.1
2	D	167	CYS	2.1
2	D	180	ASN	2.0
1	C	7	PRO	2.0
1	C	92	GLY	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.