



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 04:14 AM EDT

PDB ID : 7LKI / pdb_00007lki
Title : The crystal structure of Epitope III of HCV envelop protein E2 in complex with antibody 1H8
Authors : Deng, L.; Zhang, P.
Deposited on : 2021-02-02
Resolution : 2.00 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:

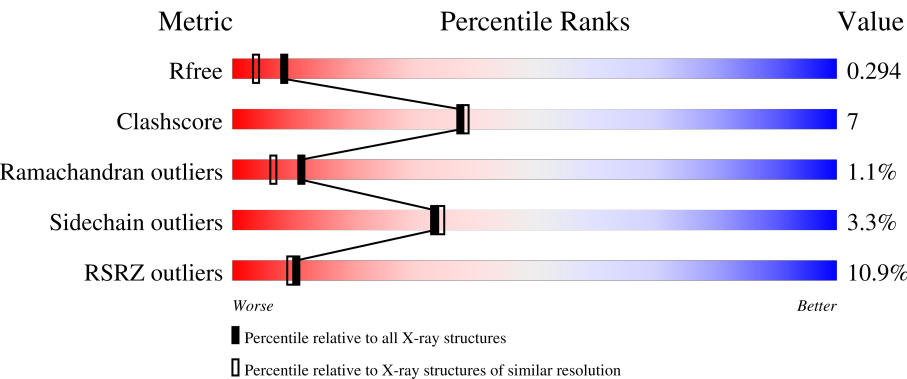
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	AAA	217	4% 84% 12% • •
1	DDD	217	16% 81% 13% • 5%
2	BBB	218	3% 82% 16% •
2	EEE	218	20% 76% 17% • •
3	CCC	15	7% 53% 47%

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Mol	Chain	Length	Quality of chain
3	FFF	15	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6986 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 antibody 1H8 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	210	Total	C	N	O	S	0	0	0
			1585	999	260	318	8			
1	DDD	206	Total	C	N	O	S	0	0	0
			1554	981	255	310	8			

- Molecule 2 is a protein called antibody 1H8 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	BBB	218	Total	C	N	O	S	0	0	0
			1686	1060	284	336	6			
2	EEE	209	Total	C	N	O	S	0	0	0
			1614	1021	268	319	6			

- Molecule 3 is a protein called Epitope III peptide GLY-ALA-PRO-THR-TYR-SER-TRP-GLY.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	CCC	8	Total	C	N	O	0	0	0
			59	39	9	11			
3	FFF	8	Total	C	N	O	0	0	0
			59	39	9	11			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	AAA	115	Total	O	0	0
			115	115		
4	BBB	116	Total	O	0	0
			116	116		
4	CCC	1	Total	O	0	0
			1	1		
4	DDD	114	Total	O	0	0
			114	114		

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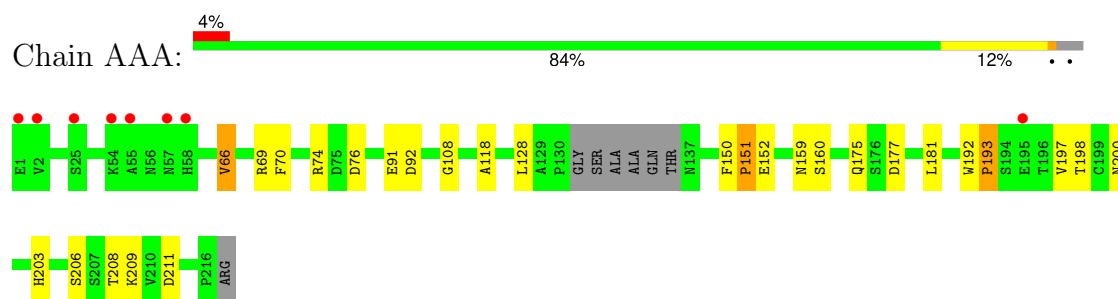
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	EEE	83	Total	O	0	0
			83	83		

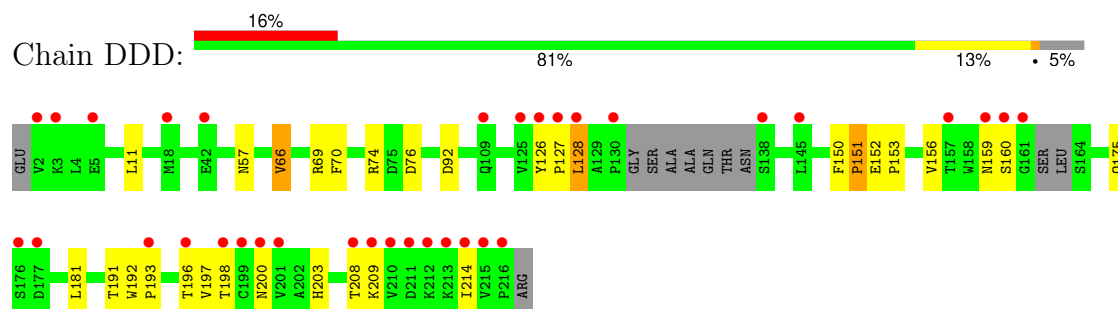
3 Residue-property plots [i](#)

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.

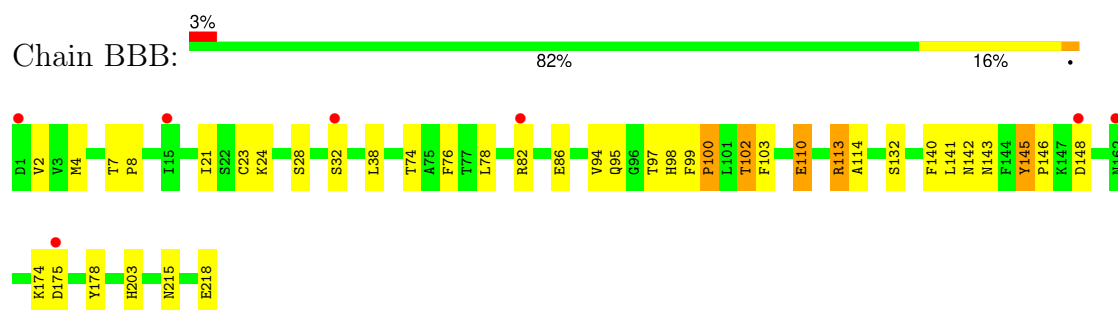
- Molecule 1: antibody 1H8 heavy chain



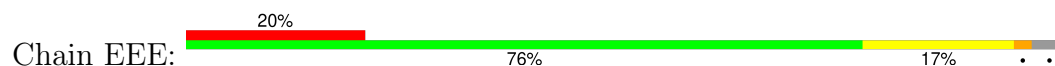
- Molecule 1: antibody 1H8 heavy chain

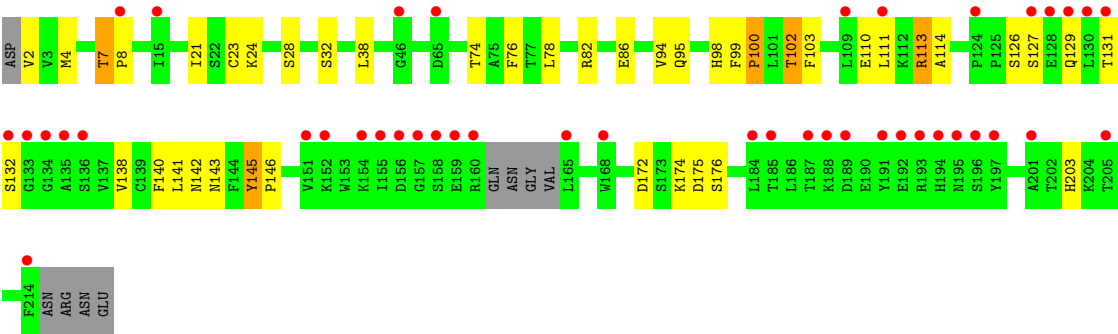


- Molecule 2: antibody 1H8 light chain

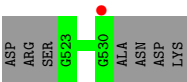


- Molecule 2: antibody 1H8 light chain

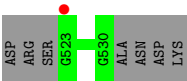




● Molecule 3: Epitope III peptide GLY-ALA-PRO-THR-TYR-SER-TRP-GLY



● Molecule 3: Epitope III peptide GLY-ALA-PRO-THR-TYR-SER-TRP-GLY



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	201.68Å 42.79Å 110.82Å 90.00° 109.53° 90.00°	Depositor
Resolution (Å)	35.70 – 2.00 35.70 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.4 (35.70-2.00) 98.4 (35.70-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.80 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.245 , 0.294 0.247 , 0.294	Depositor DCC
R_{free} test set	3030 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	22.7	Xtriage
Anisotropy	0.377	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 35.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6986	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.85% 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	AAA	1.06	0/1624	1.31	2/2215 (0.1%)
1	DDD	1.03	0/1592	1.32	0/2170
2	BBB	1.01	0/1723	1.28	1/2339 (0.0%)
2	EEE	1.03	0/1650	1.29	2/2239 (0.1%)
3	CCC	0.92	0/62	1.10	0/85
3	FFF	0.95	0/62	1.03	0/85
All	All	1.03	0/6713	1.30	5/9133 (0.1%)

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	AAA	0	5
1	DDD	0	6
2	BBB	0	4
2	EEE	0	4
All	All	0	19

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	BBB	100	PRO	N-CA-C	6.04	124.91	112.47
2	EEE	7	THR	CB-CA-C	5.49	117.50	109.48
1	AAA	108	GLY	CA-C-N	5.32	128.40	120.31
1	AAA	108	GLY	C-N-CA	5.32	128.40	120.31
2	EEE	100	PRO	N-CA-C	5.06	122.90	112.47

There are no chirality outliers.

All (19) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	AAA	150	PHE	Peptide,Mainchain
1	AAA	152	GLU	Peptide,Mainchain
1	AAA	192	TRP	Peptide
2	BBB	145	TYR	Peptide
2	BBB	7	THR	Peptide
2	BBB	99	PHE	Peptide,Mainchain
1	DDD	150	PHE	Peptide,Mainchain
1	DDD	152	GLU	Peptide,Mainchain
1	DDD	192	TRP	Peptide,Mainchain
2	EEE	145	TYR	Peptide
2	EEE	7	THR	Peptide
2	EEE	99	PHE	Peptide,Mainchain

5.2 Too-close contacts [i](#)

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	AAA	1585	0	1540	21	0
1	DDD	1554	0	1508	24	0
2	BBB	1686	0	1654	24	0
2	EEE	1614	0	1589	29	1
3	CCC	59	0	48	0	0
3	FFF	59	0	48	0	0
4	AAA	115	0	0	4	0
4	BBB	116	0	0	1	0
4	CCC	1	0	0	0	0
4	DDD	114	0	0	4	0
4	EEE	83	0	0	3	0
All	All	6986	0	6387	90	1

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 7.

All (90) 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:EEE:111:LEU:HD23	2:EEE:176:SER:OG	1.40	1.17

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BBB:95:GLN:HE22	2:BBB:98:HIS:H	1.11	0.98
2:EEE:95:GLN:HE22	2:EEE:98:HIS:H	1.05	0.95
2:EEE:95:GLN:NE2	2:EEE:98:HIS:H	1.63	0.94
2:BBB:95:GLN:NE2	2:BBB:98:HIS:H	1.69	0.89
2:EEE:95:GLN:HE22	2:EEE:98:HIS:N	1.78	0.81
1:AAA:160:SER:H	1:AAA:200:ASN:HD21	1.30	0.79
1:AAA:69:ARG:HD3	4:AAA:360:HOH:O	1.84	0.78
1:DDD:160:SER:H	1:DDD:200:ASN:HD21	1.32	0.76
2:EEE:111:LEU:HD23	2:EEE:176:SER:HG	1.48	0.74
1:DDD:74:ARG:HD3	1:DDD:76:ASP:OD1	1.89	0.73
2:BBB:95:GLN:HE22	2:BBB:98:HIS:N	1.86	0.73
1:AAA:177:ASP:CG	1:DDD:196:THR:HG23	2.16	0.69
2:EEE:127:SER:O	4:EEE:301:HOH:O	2.11	0.68
2:BBB:82:ARG:HG3	4:BBB:303:HOH:O	1.94	0.67
2:EEE:2:VAL:O	2:EEE:102:THR:HG21	1.95	0.66
2:BBB:215:ASN:HB3	2:BBB:218:GLU:HG2	1.79	0.65
2:BBB:2:VAL:O	2:BBB:102:THR:HG21	1.97	0.64
1:AAA:151:PRO:O	1:AAA:203:HIS:HE1	1.81	0.62
1:DDD:159:ASN:ND2	1:DDD:198:THR:H	1.98	0.61
1:AAA:74:ARG:HD3	1:AAA:76:ASP:OD1	2.01	0.60
1:DDD:160:SER:H	1:DDD:200:ASN:ND2	2.00	0.59
2:BBB:38:LEU:HD22	2:BBB:76:PHE:CG	2.37	0.59
1:DDD:57:ASN:ND2	4:DDD:301:HOH:O	2.12	0.59
2:EEE:113:ARG:HD3	2:EEE:114:ALA:O	2.02	0.59
1:AAA:66:VAL:HG13	1:AAA:70:PHE:HB2	1.85	0.58
2:EEE:146:PRO:HD2	2:EEE:203:HIS:CE1	2.39	0.58
1:AAA:160:SER:H	1:AAA:200:ASN:ND2	2.00	0.58
1:DDD:66:VAL:HG13	1:DDD:70:PHE:HB2	1.85	0.58
1:AAA:69:ARG:CD	4:AAA:360:HOH:O	2.45	0.58
1:DDD:151:PRO:O	1:DDD:203:HIS:HE1	1.85	0.57
1:AAA:177:ASP:OD2	1:DDD:196:THR:CG2	2.52	0.57
1:AAA:159:ASN:ND2	1:AAA:198:THR:H	2.03	0.57
1:DDD:127:PRO:O	2:EEE:126:SER:HB3	2.06	0.56
2:EEE:38:LEU:HD22	2:EEE:76:PHE:CG	2.40	0.56
2:EEE:82:ARG:NH1	4:EEE:302:HOH:O	2.37	0.55
2:BBB:28:SER:O	2:BBB:98:HIS:HE1	1.89	0.55
2:EEE:28:SER:O	2:EEE:98:HIS:HE1	1.89	0.55
1:AAA:177:ASP:OD2	1:DDD:196:THR:HG23	2.08	0.54
2:BBB:145:TYR:C	2:BBB:145:TYR:CD1	2.86	0.54
1:AAA:69:ARG:NH2	1:AAA:92:ASP:OD2	2.35	0.52
2:BBB:113:ARG:HD3	2:BBB:114:ALA:O	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:DDD:69:ARG:HD3	4:DDD:328:HOH:O	2.08	0.52
1:AAA:91:GLU:HG3	1:DDD:191:THR:CG2	2.40	0.52
1:AAA:159:ASN:HD21	1:AAA:197:VAL:HA	1.76	0.51
2:EEE:4:MET:HE3	2:EEE:23:CYS:SG	2.50	0.51
2:EEE:141:LEU:N	2:EEE:141:LEU:HD12	2.26	0.51
2:BBB:110:GLU:CD	2:BBB:178:TYR:OH	2.54	0.51
2:EEE:24:LYS:HA	2:EEE:74:THR:O	2.11	0.50
2:EEE:145:TYR:CD1	2:EEE:145:TYR:C	2.88	0.50
1:DDD:159:ASN:HD21	1:DDD:198:THR:H	1.58	0.50
1:DDD:214:ILE:O	4:DDD:302:HOH:O	2.19	0.50
2:BBB:21:ILE:HD11	2:BBB:78:LEU:HD23	1.94	0.50
1:AAA:208:THR:HG21	4:AAA:329:HOH:O	2.12	0.49
1:AAA:211:ASP:OD2	4:AAA:301:HOH:O	2.20	0.49
2:BBB:4:MET:HE3	2:BBB:23:CYS:SG	2.53	0.48
2:BBB:141:LEU:HD12	2:BBB:141:LEU:N	2.29	0.48
1:DDD:159:ASN:HD21	1:DDD:197:VAL:HA	1.80	0.47
2:EEE:131:THR:OG1	4:EEE:301:HOH:O	2.07	0.47
2:BBB:95:GLN:NE2	2:BBB:97:THR:N	2.63	0.47
2:BBB:4:MET:HG2	2:BBB:102:THR:HG23	1.98	0.46
2:BBB:24:LYS:HA	2:BBB:74:THR:O	2.16	0.46
2:EEE:142:ASN:HB3	2:EEE:143:ASN:ND2	2.32	0.45
2:EEE:21:ILE:HD11	2:EEE:78:LEU:HD23	1.98	0.45
1:DDD:203:HIS:HB3	1:DDD:208:THR:HB	1.99	0.45
2:BBB:95:GLN:HE21	2:BBB:97:THR:H	1.63	0.45
1:AAA:118:ALA:HB2	1:AAA:177:ASP:HB3	1.98	0.45
2:BBB:95:GLN:NE2	2:BBB:97:THR:H	2.14	0.45
1:AAA:181:LEU:HD12	1:AAA:181:LEU:C	2.42	0.44
1:DDD:11:LEU:HD22	1:DDD:151:PRO:HD3	1.99	0.44
2:BBB:146:PRO:O	2:BBB:203:HIS:HE1	2.00	0.44
2:EEE:4:MET:HG2	2:EEE:102:THR:HG23	2.00	0.44
1:DDD:69:ARG:NH2	1:DDD:92:ASP:OD2	2.33	0.44
2:EEE:111:LEU:CD2	2:EEE:176:SER:OG	2.35	0.44
2:EEE:95:GLN:NE2	2:EEE:98:HIS:N	2.46	0.44
2:BBB:142:ASN:HB3	2:BBB:143:ASN:ND2	2.33	0.43
1:DDD:181:LEU:C	1:DDD:181:LEU:HD12	2.43	0.43
2:BBB:94:VAL:HG22	2:BBB:103:PHE:CD2	2.53	0.43
2:BBB:140:PHE:C	2:BBB:141:LEU:HD12	2.44	0.42
1:DDD:156:VAL:O	4:DDD:303:HOH:O	2.21	0.42
2:EEE:94:VAL:HG22	2:EEE:103:PHE:CD2	2.53	0.42
2:EEE:140:PHE:C	2:EEE:141:LEU:HD12	2.43	0.42
1:AAA:66:VAL:HG13	1:AAA:70:PHE:CG	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:203:HIS:HD2	1:AAA:206:SER:OG	2.01	0.42
2:BBB:146:PRO:HD2	2:BBB:203:HIS:CE1	2.54	0.42
1:DDD:128:LEU:HD21	2:EEE:138:VAL:HG21	2.03	0.41
1:DDD:126:TYR:CE1	2:EEE:129:GLN:HA	2.56	0.41
1:DDD:127:PRO:HG2	2:EEE:126:SER:HB2	2.04	0.40
1:AAA:203:HIS:HB3	1:AAA:208:THR:HB	2.04	0.40
2:EEE:172:ASP:OD2	2:EEE:175:ASP:HB2	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:EEE:131:THR:O	2:EEE:131:THR:O[2_556]	1.80	0.40

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	AAA	206/217 (95%)	200 (97%)	4 (2%)	2 (1%)	12	8
1	DDD	200/217 (92%)	197 (98%)	0	3 (2%)	8	4
2	BBB	216/218 (99%)	210 (97%)	4 (2%)	2 (1%)	14	9
2	EEE	205/218 (94%)	198 (97%)	5 (2%)	2 (1%)	12	8
3	CCC	6/15 (40%)	6 (100%)	0	0	100	100
3	FFF	6/15 (40%)	6 (100%)	0	0	100	100
All	All	839/900 (93%)	817 (97%)	13 (2%)	9 (1%)	11	7

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	EEE	8	PRO

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Mol	Chain	Res	Type
2	BBB	8	PRO
1	DDD	151	PRO
1	DDD	153	PRO
1	AAA	151	PRO
2	BBB	100	PRO
2	EEE	100	PRO
1	DDD	193	PRO
1	AAA	193	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	AAA	181/185 (98%)	176 (97%)	5 (3%)	38	41
1	DDD	177/185 (96%)	173 (98%)	4 (2%)	44	49
2	BBB	194/194 (100%)	185 (95%)	9 (5%)	24	22
2	EEE	186/194 (96%)	179 (96%)	7 (4%)	29	29
3	CCC	5/11 (46%)	5 (100%)	0	100	100
3	FFF	5/11 (46%)	5 (100%)	0	100	100
All	All	748/780 (96%)	723 (97%)	25 (3%)	33	34

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	66	VAL
1	AAA	128	LEU
1	AAA	175	GLN
1	AAA	193	PRO
1	AAA	209	LYS
2	BBB	32	SER
2	BBB	86	GLU
2	BBB	102	THR
2	BBB	110	GLU
2	BBB	113	ARG

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Mol	Chain	Res	Type
2	BBB	132	SER
2	BBB	148	ASP
2	BBB	174	LYS
2	BBB	175	ASP
1	DDD	66	VAL
1	DDD	128	LEU
1	DDD	175	GLN
1	DDD	209	LYS
2	EEE	32	SER
2	EEE	86	GLU
2	EEE	102	THR
2	EEE	110	GLU
2	EEE	113	ARG
2	EEE	132	SER
2	EEE	174	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

There are no ligands in this entry.

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	AAA	210/217 (96%)	0.35	8 (3%) 44 43	13, 22, 42, 52	0
1	DDD	206/217 (94%)	0.94	34 (16%) 4 4	17, 28, 47, 63	0
2	BBB	218/218 (100%)	0.50	7 (3%) 50 49	15, 28, 42, 54	0
2	EEE	209/218 (95%)	1.29	43 (20%) 2 2	22, 35, 62, 84	0
3	CCC	8/15 (53%)	1.06	1 (12%) 8 7	31, 35, 40, 43	0
3	FFF	8/15 (53%)	0.91	1 (12%) 8 7	26, 28, 33, 37	0
All	All	859/900 (95%)	0.77	94 (10%) 10 9	13, 29, 50, 84	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	DDD	177	ASP	6.6
1	DDD	210	VAL	5.9
2	EEE	158	SER	5.8
2	EEE	155	ILE	5.7
2	EEE	197	TYR	5.1
1	DDD	213	LYS	4.7
1	DDD	176	SER	4.6
2	EEE	130	LEU	4.5
2	EEE	195	ASN	4.5
2	EEE	132	SER	4.3
2	EEE	187	THR	4.2
2	EEE	131	THR	4.1
2	EEE	184	LEU	4.1
1	DDD	212	LYS	4.1
1	DDD	159	ASN	4.0
1	DDD	215	VAL	3.9
1	DDD	211	ASP	3.9
2	BBB	1	ASP	3.8
2	EEE	191	TYR	3.7

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Mol	Chain	Res	Type	RSRZ
2	EEE	185	THR	3.7
2	EEE	193	ARG	3.4
2	EEE	127	SER	3.4
1	AAA	195	GLU	3.3
2	BBB	15	ILE	3.3
1	DDD	208	THR	3.2
1	AAA	57	ASN	3.2
1	DDD	161	GLY	3.2
1	DDD	199	CYS	3.1
1	DDD	128	LEU	3.1
2	EEE	65	ASP	3.1
1	AAA	58	HIS	3.0
1	DDD	157	THR	3.0
1	DDD	214	ILE	3.0
2	EEE	188	LYS	3.0
2	EEE	135	ALA	3.0
2	EEE	46	GLY	3.0
1	DDD	216	PRO	3.0
2	EEE	111	LEU	2.9
1	DDD	127	PRO	2.9
1	AAA	54	LYS	2.9
2	EEE	196	SER	2.9
2	EEE	160	ARG	2.8
2	EEE	194	HIS	2.8
2	EEE	189	ASP	2.8
2	EEE	214	PHE	2.8
1	DDD	5	GLU	2.7
1	DDD	198	THR	2.7
2	EEE	8	PRO	2.7
2	BBB	148	ASP	2.7
1	DDD	193	PRO	2.7
2	EEE	156	ASP	2.7
2	EEE	134	GLY	2.7
2	EEE	157	GLY	2.6
1	DDD	126	TYR	2.6
1	DDD	201	VAL	2.5
2	EEE	205	THR	2.5
1	DDD	160	SER	2.5
1	DDD	145	LEU	2.5
2	BBB	162	ASN	2.5
1	DDD	209	LYS	2.5
2	EEE	15	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
2	BBB	32	SER	2.4
2	EEE	168	TRP	2.4
1	AAA	55	ALA	2.4
2	EEE	129	GLN	2.4
1	DDD	138	SER	2.3
1	AAA	2	VAL	2.3
2	EEE	152	LYS	2.3
1	DDD	42	GLU	2.2
1	DDD	109	GLN	2.2
1	DDD	2	VAL	2.2
2	EEE	159	GLU	2.2
1	DDD	125	VAL	2.2
3	FFF	523	GLY	2.2
1	DDD	3	LYS	2.2
1	DDD	200	ASN	2.2
2	BBB	175	ASP	2.2
1	DDD	196	THR	2.2
2	EEE	154	LYS	2.1
2	EEE	128	GLU	2.1
1	DDD	130	PRO	2.1
2	EEE	109	LEU	2.1
2	EEE	192	GLU	2.1
2	EEE	133	GLY	2.1
2	EEE	151	VAL	2.1
1	AAA	25	SER	2.1
2	EEE	165	LEU	2.1
1	AAA	1	GLU	2.0
2	EEE	136	SER	2.1
2	BBB	82	ARG	2.0
3	CCC	530	GLY	2.0
1	DDD	18	MET	2.0
2	EEE	201	ALA	2.0
2	EEE	124	PRO	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.