



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:28 AM UTC

PDB ID : 7NDQ / pdb_00007ndq
Title : Gag:02 TCR in complex with HLA-E.
Authors : Pengelly, R.J.; Robinson, R.A.
Deposited on : 2021-02-02
Resolution : 2.55 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

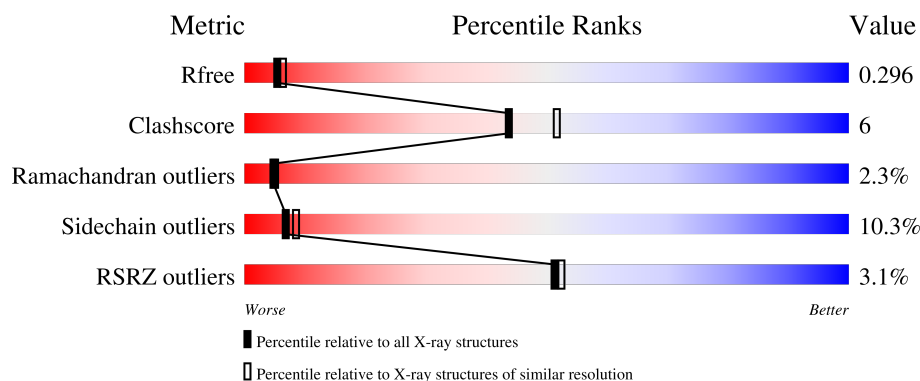
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1853 (2.58-2.54)
Clashscore	190562	1897 (2.58-2.54)
Ramachandran outliers	187476	1875 (2.58-2.54)
Sidechain outliers	187428	1875 (2.58-2.54)
RSRZ outliers	180081	1853 (2.58-2.54)

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	277	5% 77% 19% .
2	BBB	100	72% 25% .
3	CCC	9	33% 56% 11%
4	DDD	199	5% 65% 23% 5% 7% .
5	EEE	244	% 77% 21% ..

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6573 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 HLA class I histocompatibility antigen, alpha chain E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	276	Total	C	N	O	S	0	1	0
			2253	1408	403	435	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AAA	0	MET	-	initiating methionine	UNP P13747

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	BBB	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
BBB	0	MET	-	initiating methionine	UNP P61769

- Molecule 3 is a protein called Gag6V.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	CCC	9	Total	C	N	O	S	0	0	0
			74	48	12	13	1			

- Molecule 4 is a protein called TCR Gag:02-alpha,TCR Gag:02-alpha,TCR Gag:02-alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	DDD	185	Total	C	N	O	S	0	2	0
			1452	908	240	296	8			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
DDD	0	MET	-	initiating methionine	UNP A0A0B4J268
DDD	108	SER	-	linker	UNP A0A0B4J268
DDD	109	SER	-	linker	UNP A0A0B4J268
DDD	110	PHE	-	linker	UNP A0A0B4J268
DDD	129	ASN	-	linker	UNP A0A075B6U7
DDD	176	CYS	THR	engineered mutation	UNP P0DSE1

- Molecule 5 is a protein called T cell receptor beta variable 7-9,M1-specific T cell receptor beta chain,T cell receptor beta variable 7-9,M1-specific T cell receptor beta chain,T cell receptor beta variable 7-9,M1-specific T cell receptor beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	EEE	241	Total	C	N	O	S	0	0	0
			1944	1219	347	372	6			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
EEE	0	MET	-	initiating methionine	UNP P04435
EEE	109	GLY	-	linker	UNP P04435
EEE	110	ARG	-	linker	UNP P04435
EEE	113	GLU	-	linker	UNP P04435
EEE	132	ASN	LYS	engineered mutation	UNP K7N5M4
EEE	133	LYS	ASN	engineered mutation	UNP K7N5M4
EEE	217	ASP	ASN	engineered mutation	UNP K7N5M4

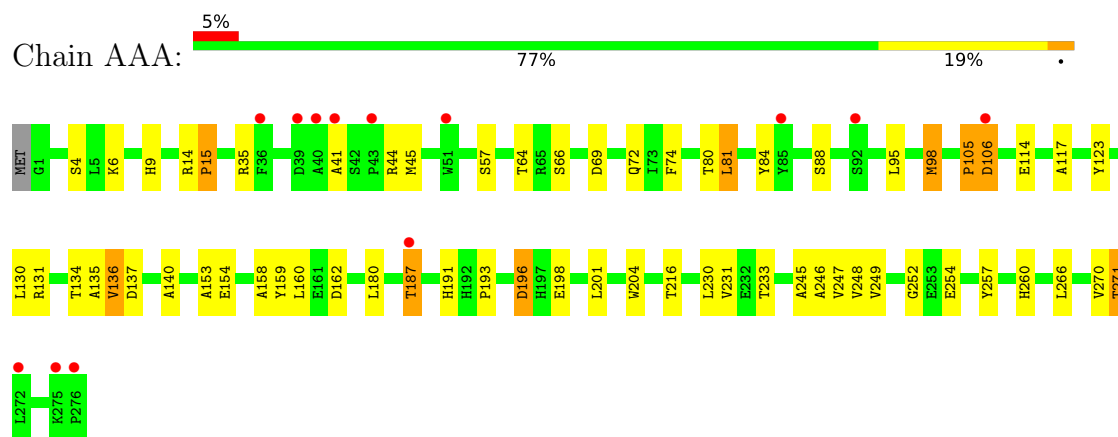
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	AAA	8	Total	O	0	0
			8	8		
6	BBB	3	Total	O	0	0
			3	3		
6	DDD	1	Total	O	0	0
			1	1		
6	EEE	1	Total	O	0	0
			1	1		

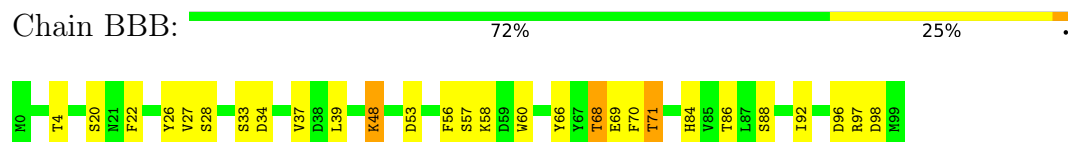
3 Residue-property plots

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: HLA class I histocompatibility antigen, alpha chain E



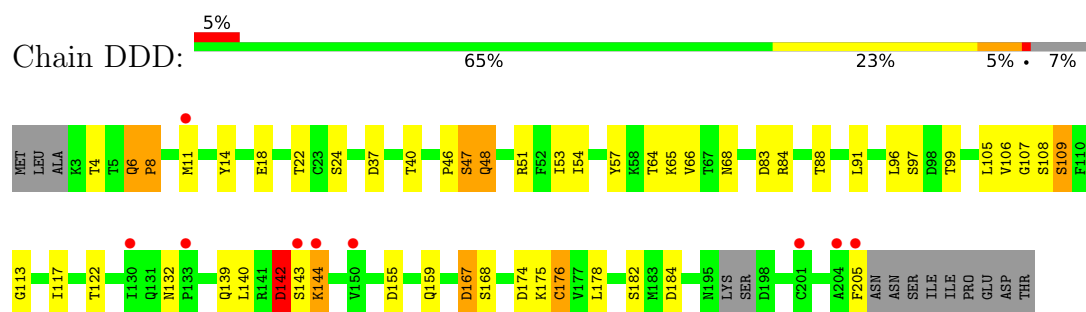
- Molecule 2: Beta-2-microglobulin



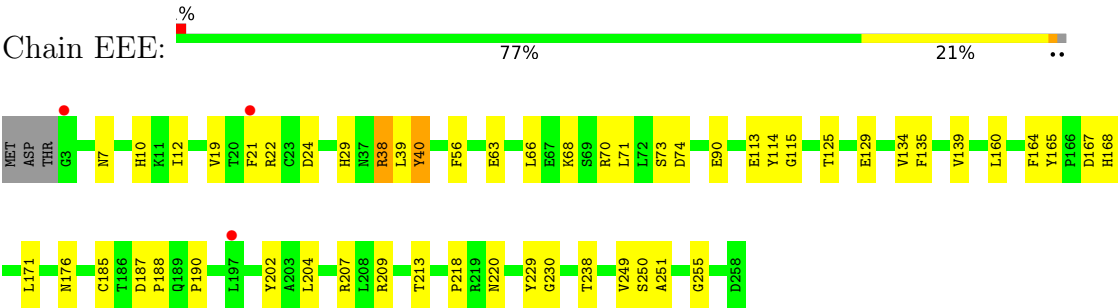
- Molecule 3: Gag6V



- Molecule 4: TCR Gag:02-alpha, TCR Gag:02-alpha, TCR Gag:02-alpha



● Molecule 5: T cell receptor beta variable 7-9,M1-specific T cell receptor beta chain,T cell receptor beta variable 7-9,M1-specific T cell receptor beta chain,T cell receptor beta variable 7-9,M1-specific T cell receptor beta chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	89.39Å 89.39Å 293.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	76.44 – 2.55 76.44 – 2.55	Depositor EDS
% Data completeness (in resolution range)	100.0 (76.44-2.55) 99.8 (76.44-2.55)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.98 (at 2.55Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.240 , 0.296 0.246 , 0.296	Depositor DCC
R_{free} test set	1952 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	78.1	Xtriage
Anisotropy	0.451	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 47.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6573	wwPDB-VP
Average B, all atoms (Å ²)	97.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.57% 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	1.04	0/2320	1.38	1/3154 (0.0%)
2	BBB	1.00	0/860	1.34	3/1162 (0.3%)
3	CCC	1.00	0/75	1.46	1/99 (1.0%)
4	DDD	1.11	0/1487	1.46	4/2017 (0.2%)
5	EEE	1.05	1/1995 (0.1%)	1.37	1/2708 (0.0%)
All	All	1.05	1/6737 (0.0%)	1.39	10/9140 (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	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	EEE	10	HIS	CE1-NE2	5.06	1.37	1.32

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	DDD	8	PRO	CA-N-CD	-10.21	97.71	112.00
4	DDD	6	GLN	C-N-CD	-6.88	96.81	125.00
4	DDD	142	ASP	CA-CB-CG	6.28	118.88	112.60
1	AAA	233	THR	CA-C-O	-5.48	115.53	121.44
4	DDD	167	ASP	CA-CB-CG	5.47	118.08	112.60
5	EEE	7	ASN	CB-CA-C	-5.47	101.82	110.02
2	BBB	26	TYR	CA-C-N	5.34	128.24	120.98
2	BBB	26	TYR	C-N-CA	5.34	128.24	120.98
2	BBB	71	THR	CB-CA-C	5.24	117.15	110.34
3	CCC	3	TYR	CB-CA-C	5.01	118.42	109.65

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	AAA	136	VAL	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	AAA	2253	0	2092	30	0
2	BBB	837	0	803	12	0
3	CCC	74	0	81	5	0
4	DDD	1452	0	1387	24	0
5	EEE	1944	0	1846	24	0
6	AAA	8	0	0	0	0
6	BBB	3	0	0	0	0
6	DDD	1	0	0	0	0
6	EEE	1	0	0	0	0
All	All	6573	0	6209	78	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 (78) 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:BBB:4:THR:HA	2:BBB:86:THR:HG21	1.54	0.87
4:DDD:8:PRO:O	4:DDD:122:THR:OG1	2.07	0.72
4:DDD:37:ASP:O	4:DDD:84:ARG:NH2	2.21	0.72
2:BBB:39:LEU:HD12	2:BBB:68:THR:HG22	1.70	0.72
1:AAA:260:HIS:ND1	1:AAA:271:THR:HG22	2.07	0.69
1:AAA:44:ARG:HA	1:AAA:64:THR:HG23	1.74	0.69
1:AAA:260:HIS:ND1	1:AAA:271:THR:CG2	2.60	0.65
4:DDD:139:GLN:C	4:DDD:140:LEU:HD12	2.24	0.62
1:AAA:98:MET:HE3	2:BBB:60:TRP:CZ3	2.36	0.61
4:DDD:22:THR:HG22	4:DDD:88:THR:HG23	1.82	0.61
2:BBB:4:THR:HA	2:BBB:86:THR:CG2	2.32	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:74:PHE:CD2	1:AAA:95:LEU:HD23	2.39	0.57
1:AAA:114:GLU:OE2	3:CCC:3:TYR:OH	2.19	0.56
4:DDD:113:GLY:O	5:EEE:38:ARG:NH1	2.37	0.56
1:AAA:106:ASP:OD1	1:AAA:106:ASP:N	2.36	0.55
1:AAA:136:VAL:HG23	1:AAA:137:ASP:HB2	1.89	0.55
4:DDD:111:ASN:N	4:DDD:111:ASN:OD1	2.40	0.55
4:DDD:109:SER:OG	5:EEE:113:GLU:HG3	2.06	0.55
1:AAA:117:ALA:HB2	2:BBB:60:TRP:CE2	2.42	0.55
4:DDD:108:SER:O	4:DDD:112:GLN:HB2	2.07	0.55
1:AAA:252:GLY:N	1:AAA:254:GLU:OE2	2.39	0.54
4:DDD:48:GLN:HB3	4:DDD:51:ARG:HH21	1.73	0.54
5:EEE:39:LEU:C	5:EEE:39:LEU:HD23	2.32	0.54
4:DDD:107:GLY:HA3	4:DDD:112:GLN:NE2	2.24	0.53
3:CCC:6:VAL:O	5:EEE:114:TYR:HB2	2.09	0.53
5:EEE:29:HIS:HE1	5:EEE:115:GLY:O	1.91	0.53
4:DDD:155:ASP:OD1	5:EEE:209:ARG:NH1	2.42	0.53
5:EEE:139:VAL:CG2	5:EEE:249:VAL:HG12	2.40	0.52
4:DDD:53:ILE:O	4:DDD:68[B]:ASN:ND2	2.43	0.51
1:AAA:249:VAL:HG12	1:AAA:257:TYR:CZ	2.45	0.51
5:EEE:167:ASP:O	5:EEE:167:ASP:CG	2.50	0.51
4:DDD:178:LEU:HB3	5:EEE:185:CYS:HB2	1.94	0.50
1:AAA:135:ALA:HB1	1:AAA:140:ALA:CB	2.41	0.49
5:EEE:38:ARG:HA	5:EEE:56:PHE:O	2.11	0.49
5:EEE:171:LEU:C	5:EEE:171:LEU:HD23	2.37	0.49
1:AAA:153:ALA:HA	4:DDD:57:TYR:HB2	1.94	0.49
4:DDD:14:TYR:OH	4:DDD:159:GLN:NE2	2.46	0.48
4:DDD:96:LEU:O	4:DDD:99:THR:OG1	2.22	0.48
2:BBB:39:LEU:CD1	2:BBB:68:THR:HG22	2.43	0.48
5:EEE:218:PRO:HA	5:EEE:255:GLY:O	2.13	0.48
1:AAA:4:SER:OG	1:AAA:6:LYS:HE2	2.14	0.48
5:EEE:168:HIS:HB3	5:EEE:229:TYR:HB2	1.96	0.47
1:AAA:230:LEU:HD13	1:AAA:245:ALA:HB2	1.96	0.47
1:AAA:35:ARG:HD3	2:BBB:53:ASP:CG	2.40	0.47
1:AAA:196:ASP:OD1	1:AAA:196:ASP:N	2.47	0.47
1:AAA:201:LEU:O	1:AAA:246:ALA:HA	2.16	0.46
5:EEE:135:PHE:O	5:EEE:164:PHE:HA	2.16	0.46
4:DDD:18:GLU:HA	4:DDD:91:LEU:O	2.16	0.46
1:AAA:266:LEU:HD13	1:AAA:270:VAL:HG13	1.97	0.45
1:AAA:69:ASP:OD2	5:EEE:38:ARG:NH2	2.49	0.45
4:DDD:4:THR:HG21	4:DDD:106:VAL:HB	1.98	0.45
5:EEE:187:ASP:HB2	5:EEE:204:LEU:HD12	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:DDD:105:LEU:HD23	4:DDD:117:ILE:O	2.18	0.44
1:AAA:80:THR:HG22	1:AAA:84:TYR:CZ	2.53	0.44
1:AAA:191:HIS:CD2	1:AAA:193:PRO:HD3	2.53	0.43
5:EEE:176:ASN:HD21	5:EEE:220:ASN:HA	1.82	0.43
2:BBB:37:VAL:HG11	2:BBB:66:TYR:CD1	2.53	0.43
2:BBB:96:ASP:O	2:BBB:98:ASP:N	2.51	0.43
1:AAA:187:THR:HA	1:AAA:204:TRP:O	2.19	0.43
5:EEE:40:TYR:N	5:EEE:40:TYR:CD1	2.86	0.43
1:AAA:131:ARG:NH1	1:AAA:154:GLU:OE2	2.52	0.42
5:EEE:190:PRO:HB2	5:EEE:202:TYR:HD1	1.84	0.42
1:AAA:159:TYR:OH	3:CCC:1:ARG:O	2.35	0.42
2:BBB:22:PHE:CE1	2:BBB:69:GLU:HB3	2.53	0.42
1:AAA:9:HIS:CE1	3:CCC:2:MET:HE3	2.55	0.42
5:EEE:139:VAL:HG23	5:EEE:249:VAL:HG12	2.01	0.42
4:DDD:46:PRO:O	4:DDD:47:SER:OG	2.27	0.42
1:AAA:84:TYR:OH	3:CCC:9:LEU:O	2.34	0.42
4:DDD:176:CYS:C	5:EEE:185:CYS:SG	3.03	0.42
1:AAA:81:LEU:HD11	1:AAA:123:TYR:CZ	2.55	0.41
1:AAA:98:MET:HE2	2:BBB:56:PHE:HE1	1.85	0.41
1:AAA:158:ALA:O	1:AAA:162:ASP:HB2	2.20	0.41
4:DDD:144:LYS:NZ	5:EEE:251:ALA:HB1	2.35	0.41
2:BBB:84:HIS:ND1	2:BBB:86:THR:HG22	2.36	0.41
4:DDD:140:LEU:HD12	4:DDD:140:LEU:N	2.36	0.41
5:EEE:134:VAL:HG13	5:EEE:165:TYR:O	2.21	0.41
4:DDD:14:TYR:OH	4:DDD:159:GLN:CD	2.64	0.40
5:EEE:12:ILE:HD12	5:EEE:230:GLY:HA2	2.04	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	AAA	275/277 (99%)	242 (88%)	29 (10%)	4 (2%)	8	10
2	BBB	98/100 (98%)	85 (87%)	8 (8%)	5 (5%)	1	1
3	CCC	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
4	DDD	183/199 (92%)	151 (82%)	25 (14%)	7 (4%)	2	1
5	EEE	239/244 (98%)	216 (90%)	21 (9%)	2 (1%)	16	22
All	All	802/829 (97%)	700 (87%)	84 (10%)	18 (2%)	5	5

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	BBB	33	SER
4	DDD	143	SER
1	AAA	105	PRO
2	BBB	20	SER
2	BBB	58	LYS
4	DDD	47	SER
4	DDD	184	ASP
2	BBB	97	ARG
4	DDD	6	GLN
4	DDD	142	ASP
1	AAA	160	LEU
2	BBB	48	LYS
5	EEE	66	LEU
1	AAA	15	PRO
4	DDD	168	SER
4	DDD	176	CYS
1	AAA	41	ALA
5	EEE	188	PRO

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	AAA	238/238 (100%)	215 (90%)	23 (10%)	8	10

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	BBB	95/95 (100%)	85 (90%)	10 (10%)	6	8
3	CCC	9/9 (100%)	7 (78%)	2 (22%)	1	1
4	DDD	169/180 (94%)	148 (88%)	21 (12%)	4	4
5	EEE	211/214 (99%)	191 (90%)	20 (10%)	8	10
All	All	722/736 (98%)	646 (90%)	76 (10%)	7	8

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	14	ARG
1	AAA	15	PRO
1	AAA	45	MET
1	AAA	57	SER
1	AAA	66	SER
1	AAA	72	GLN
1	AAA	81	LEU
1	AAA	88[A]	SER
1	AAA	88[B]	SER
1	AAA	98	MET
1	AAA	105	PRO
1	AAA	106	ASP
1	AAA	130	LEU
1	AAA	134	THR
1	AAA	180	LEU
1	AAA	187	THR
1	AAA	196	ASP
1	AAA	198	GLU
1	AAA	216	THR
1	AAA	231	VAL
1	AAA	247	VAL
1	AAA	248	VAL
1	AAA	271	THR
2	BBB	27	VAL
2	BBB	28	SER
2	BBB	34	ASP
2	BBB	48	LYS
2	BBB	57	SER
2	BBB	68	THR
2	BBB	70	PHE
2	BBB	71	THR
2	BBB	88	SER

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Mol	Chain	Res	Type
2	BBB	92	ILE
3	CCC	3	TYR
3	CCC	4	SER
4	DDD	11	MET
4	DDD	24	SER
4	DDD	40	THR
4	DDD	48	GLN
4	DDD	54	ILE
4	DDD	64	THR
4	DDD	65	LYS
4	DDD	66	VAL
4	DDD	83[A]	ASP
4	DDD	83[B]	ASP
4	DDD	97	SER
4	DDD	109	SER
4	DDD	111	ASN
4	DDD	132	ASN
4	DDD	142	ASP
4	DDD	144	LYS
4	DDD	167	ASP
4	DDD	174	ASP
4	DDD	175	LYS
4	DDD	182	SER
4	DDD	205	PHE
5	EEE	19	VAL
5	EEE	21	PHE
5	EEE	22	ARG
5	EEE	24	ASP
5	EEE	38	ARG
5	EEE	40	TYR
5	EEE	63	GLU
5	EEE	68	LYS
5	EEE	70	ARG
5	EEE	71	LEU
5	EEE	73	SER
5	EEE	74	ASP
5	EEE	90	GLU
5	EEE	125	THR
5	EEE	129	GLU
5	EEE	160	LEU
5	EEE	207	ARG
5	EEE	213	THR

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Mol	Chain	Res	Type
5	EEE	238	THR
5	EEE	250	SER

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

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	AAA	276/277 (99%)	0.47	13 (4%) 36 37	46, 86, 117, 134	1 (0%)
2	BBB	100/100 (100%)	0.30	0 100 100	64, 86, 103, 119	0
3	CCC	9/9 (100%)	0.36	0 100 100	71, 77, 87, 88	0
4	DDD	185/199 (92%)	0.59	9 (4%) 35 35	72, 118, 150, 172	2 (1%)
5	EEE	241/244 (98%)	0.21	3 (1%) 76 78	67, 94, 133, 144	0
All	All	811/829 (97%)	0.40	25 (3%) 51 52	46, 93, 140, 172	3 (0%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	DDD	205	PHE	4.0
4	DDD	143	SER	3.9
4	DDD	204	ALA	3.7
4	DDD	144	LYS	3.4
1	AAA	272	LEU	3.0
1	AAA	276	PRO	2.7
1	AAA	36	PHE	2.6
1	AAA	43	PRO	2.6
1	AAA	51	TRP	2.6
5	EEE	197	LEU	2.6
4	DDD	130	ILE	2.6
1	AAA	275	LYS	2.5
1	AAA	41	ALA	2.4
4	DDD	133	PRO	2.3
1	AAA	40	ALA	2.3
1	AAA	39	ASP	2.3
4	DDD	150	VAL	2.3
4	DDD	11	MET	2.2
1	AAA	85	TYR	2.2
5	EEE	3	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
4	DDD	201	CYS	2.1
5	EEE	21	PHE	2.1
1	AAA	106	ASP	2.1
1	AAA	92	SER	2.1
1	AAA	187	THR	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.