



# Full wwPDB X-ray Structure Validation Report ⓘ

May 20, 2026 – 08:12 PM EDT

PDB ID : 9ZMA / pdb\_00009zma  
Title : Crystal structure of Fab 7160 in complex with major repeat region (NANP6) from circumsporozoite protein  
Authors : Jain, M.; Wilson, I.A.  
Deposited on : 2025-12-09  
Resolution : 2.33 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](mailto: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>.

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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

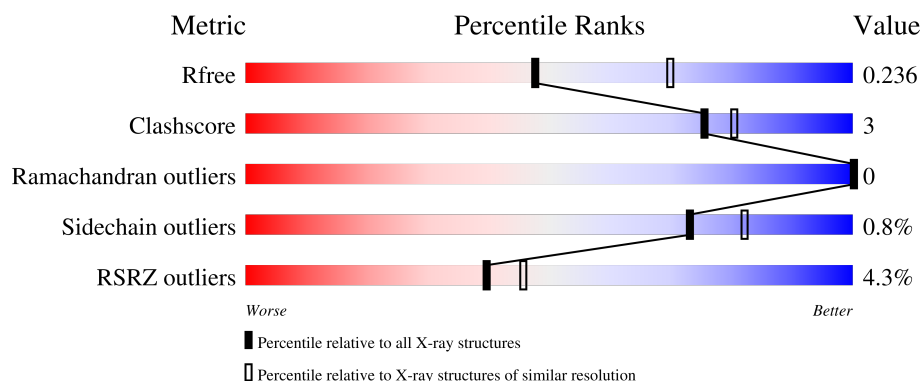
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

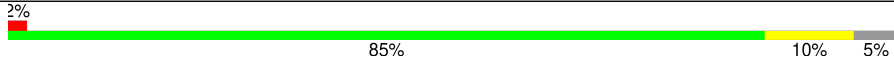


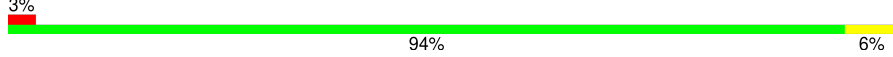
The reported resolution of this entry is 2.33 Å.

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.



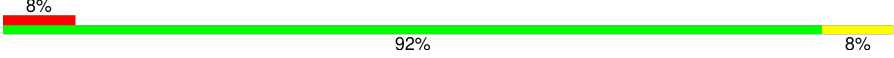

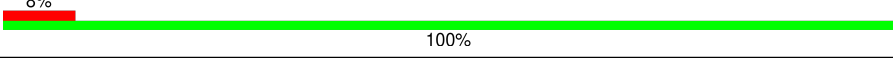
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	3031 (2.36-2.32)
Clashscore	190562	3127 (2.36-2.32)
Ramachandran outliers	187476	3095 (2.36-2.32)
Sidechain outliers	187428	3095 (2.36-2.32)
RSRZ outliers	180081	3033 (2.36-2.32)

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  $\geq 3$ , 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	230	
1	C	230	
1	E	230	
1	G	230	
2	B	219	

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Mol	Chain	Length	Quality of chain
2	D	219	
2	F	219	
2	H	219	
3	I	24	
3	J	24	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13936 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 Heavy Chain of Fab 7160.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	220	Total	C	N	O	S	0	0	0
			1640	1038	275	321	6			
1	C	219	Total	C	N	O	S	0	0	0
			1631	1032	273	320	6			
1	E	217	Total	C	N	O	S	0	0	0
			1623	1029	272	316	6			
1	G	218	Total	C	N	O	S	0	0	0
			1627	1031	273	317	6			

- Molecule 2 is a protein called Light Chain of Fab 7160.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	218	Total	C	N	O	S	0	0	0
			1697	1067	284	340	6			
2	D	219	Total	C	N	O	S	0	0	0
			1703	1070	285	341	7			
2	F	219	Total	C	N	O	S	0	0	0
			1703	1070	285	341	7			
2	H	219	Total	C	N	O	S	0	0	0
			1703	1070	285	341	7			

- Molecule 3 is a protein called Circumsporozoite protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	I	21	Total	C	N	O	0	0	0
			148	84	32	32			
3	J	24	Total	C	N	O	0	0	0
			168	96	36	36			

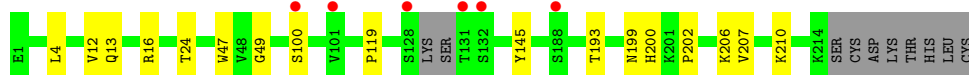
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	51	Total 51	O 51	0	0
4	B	45	Total 45	O 45	0	0
4	C	50	Total 50	O 50	0	0
4	D	29	Total 29	O 29	0	0
4	E	28	Total 28	O 28	0	0
4	F	21	Total 21	O 21	0	0
4	G	41	Total 41	O 41	0	0
4	H	16	Total 16	O 16	0	0
4	I	4	Total 4	O 4	0	0
4	J	8	Total 8	O 8	0	0

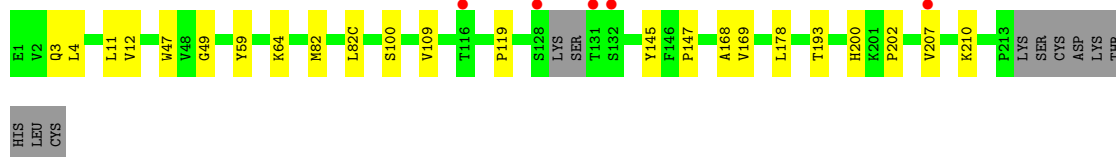
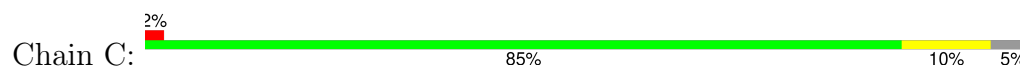
### 3 Residue-property plots [i](#)

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.

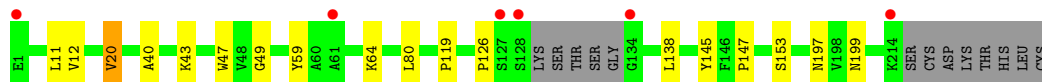
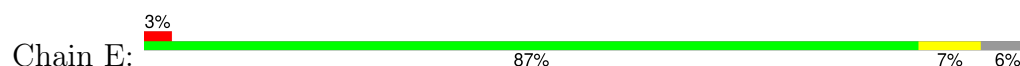
- Molecule 1: Heavy Chain of Fab 7160



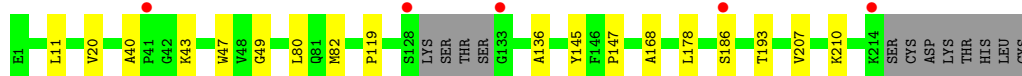
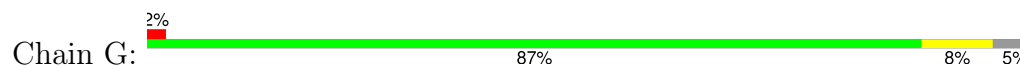
- Molecule 1: Heavy Chain of Fab 7160



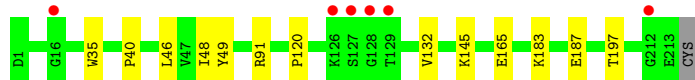
- Molecule 1: Heavy Chain of Fab 7160



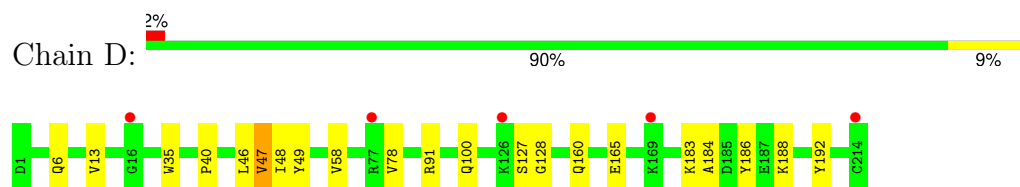
- Molecule 1: Heavy Chain of Fab 7160



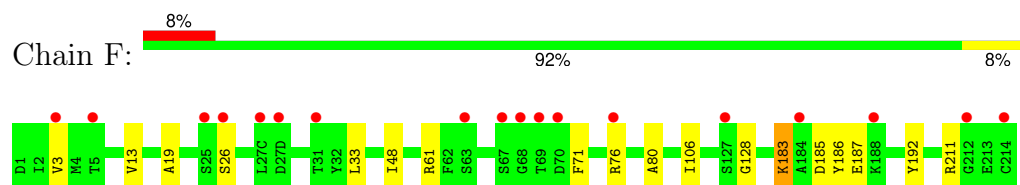
- Molecule 2: Light Chain of Fab 7160



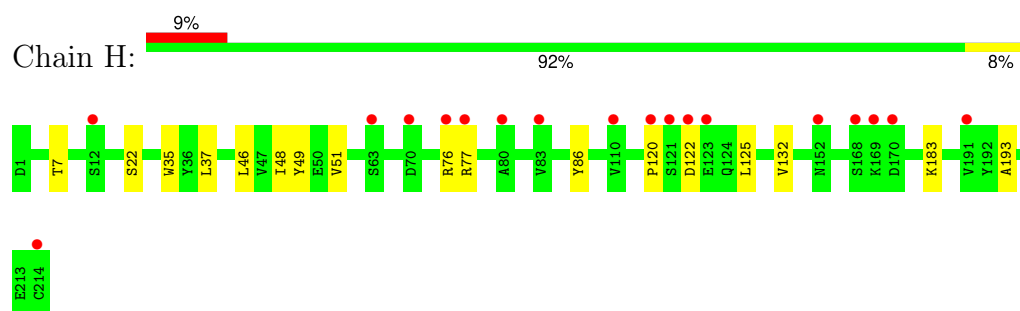
- Molecule 2: Light Chain of Fab 7160



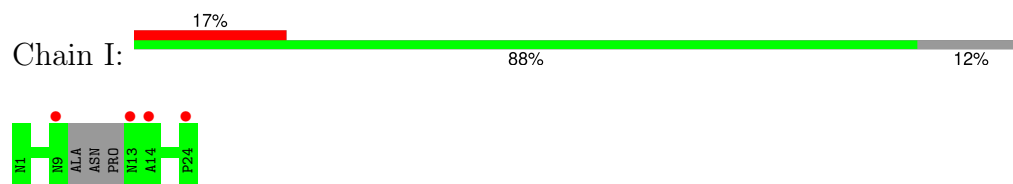
- Molecule 2: Light Chain of Fab 7160



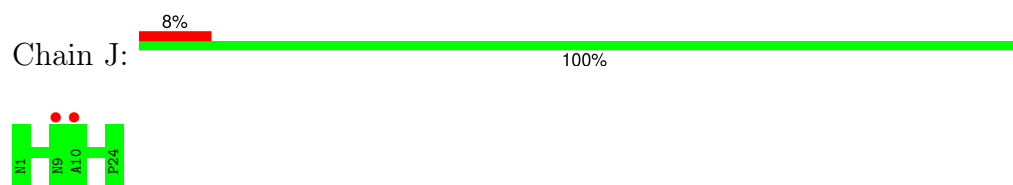
- Molecule 2: Light Chain of Fab 7160



- Molecule 3: Circumsporozoite protein



- Molecule 3: Circumsporozoite protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	103.21Å 73.13Å 134.37Å 90.00° 102.25° 90.00°	Depositor
Resolution (Å)	41.74 – 2.33 41.74 – 2.33	Depositor EDS
% Data completeness (in resolution range)	97.6 (41.74-2.33) 96.1 (41.74-2.33)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.48 (at 2.34Å)	Xtriage
Refinement program	PHENIX (1.21.2_5419: ???)	Depositor
R, $R_{free}$	0.207 , 0.236 0.207 , 0.236	Depositor DCC
$R_{free}$ test set	2000 reflections (2.44%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.9	Xtriage
Anisotropy	0.503	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 26.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	13936	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

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	0.12	0/1679	0.30	0/2285
1	C	0.13	0/1670	0.30	0/2274
1	E	0.12	0/1662	0.28	0/2262
1	G	0.13	0/1666	0.31	0/2267
2	B	0.12	0/1734	0.30	0/2354
2	D	0.12	0/1740	0.30	0/2362
2	F	0.12	0/1740	0.30	0/2362
2	H	0.15	0/1740	0.35	0/2362
3	I	0.11	0/151	0.27	0/210
3	J	0.12	0/173	0.30	0/243
All	All	0.13	0/13955	0.31	0/18981

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	A	1640	0	1618	9	0
1	C	1631	0	1605	13	0
1	E	1623	0	1603	9	0
1	G	1627	0	1606	9	0
2	B	1697	0	1659	7	0
2	D	1703	0	1664	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	1703	0	1664	10	0
2	H	1703	0	1664	11	0
3	I	148	0	127	0	0
3	J	168	0	146	0	0
4	A	51	0	0	0	0
4	B	45	0	0	0	0
4	C	50	0	0	0	0
4	D	29	0	0	0	0
4	E	28	0	0	0	0
4	F	21	0	0	0	0
4	G	41	0	0	0	0
4	H	16	0	0	0	0
4	I	4	0	0	0	0
4	J	8	0	0	0	0
All	All	13936	0	13356	80	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 3.

All (80) 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:E:40:ALA:H	1:E:43:LYS:HE2	1.42	0.85
2:D:47:VAL:HG22	2:D:48:ILE:HG12	1.71	0.71
2:F:3:VAL:HG22	2:F:26:SER:HB3	1.72	0.71
1:E:11:LEU:HB2	1:E:147:PRO:HG3	1.77	0.64
1:G:11:LEU:HB2	1:G:147:PRO:HG3	1.79	0.64
2:F:61:ARG:HA	2:F:76:ARG:HD3	1.78	0.63
2:B:46:LEU:HD21	2:B:49:TYR:HB3	1.80	0.62
2:H:120:PRO:HD3	2:H:132:VAL:HG22	1.82	0.61
2:D:6:GLN:H	2:D:100:GLN:NE2	1.97	0.60
1:C:119:PRO:HB3	1:C:145:TYR:HB3	1.85	0.59
1:E:20:VAL:HG13	1:E:80:LEU:HB3	1.84	0.59
1:E:59:TYR:HB2	1:E:64:LYS:HG3	1.84	0.59
2:F:186:TYR:HA	2:F:192:TYR:OH	2.04	0.58
1:E:119:PRO:HB3	1:E:145:TYR:HB3	1.86	0.57
2:D:6:GLN:H	2:D:100:GLN:HE22	1.53	0.56
2:D:46:LEU:HD21	2:D:49:TYR:HB3	1.88	0.55
1:G:20:VAL:HG13	1:G:80:LEU:HB3	1.87	0.55
1:A:199:ASN:HD22	1:A:206:LYS:HG2	1.71	0.55
1:C:82:MET:HE1	1:C:109:VAL:HG21	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:80:ALA:HA	2:F:106:ILE:HG13	1.89	0.53
1:E:47:TRP:CZ2	1:E:49:GLY:HA2	2.43	0.53
2:D:91:ARG:O	2:D:91:ARG:HG2	2.07	0.52
2:H:122:ASP:HA	2:H:125:LEU:HG	1.91	0.52
2:D:47:VAL:HG23	2:D:58:VAL:HG11	1.92	0.51
2:F:13:VAL:HG21	2:F:19:ALA:HB2	1.93	0.51
1:G:193:THR:HG23	1:G:210:LYS:HE3	1.93	0.51
1:A:47:TRP:CZ2	1:A:49:GLY:HA2	2.46	0.51
2:H:193:ALA:HB2	2:H:208:SER:HB3	1.93	0.51
1:A:200:HIS:CD2	1:A:202:PRO:HD2	2.45	0.50
1:G:40:ALA:HB3	1:G:43:LYS:HB2	1.94	0.50
1:A:119:PRO:HB3	1:A:145:TYR:HB3	1.93	0.50
1:C:100:SER:O	2:D:91:ARG:HD2	2.11	0.50
2:H:76:ARG:CZ	2:H:77:ARG:NH2	2.75	0.50
2:B:35:TRP:HB2	2:B:48:ILE:HB	1.94	0.49
2:F:61:ARG:HA	2:F:76:ARG:HH11	1.77	0.49
1:C:200:HIS:CD2	1:C:202:PRO:HD2	2.48	0.48
1:C:12:VAL:HG11	1:C:82(C):LEU:HD13	1.96	0.47
1:A:13:GLN:HB2	1:A:16:ARG:HG3	1.96	0.46
1:C:47:TRP:CZ2	1:C:49:GLY:HA2	2.51	0.46
2:D:40:PRO:HB3	2:D:165:GLU:HG3	1.97	0.46
1:G:119:PRO:HB3	1:G:145:TYR:HB3	1.98	0.46
1:C:3:GLN:O	1:C:4:LEU:HD23	2.16	0.46
2:F:185:ASP:C	2:F:187:GLU:H	2.24	0.46
1:C:82:MET:HB3	1:C:82(C):LEU:HD21	1.97	0.46
1:E:40:ALA:HB3	1:E:43:LYS:HG2	1.97	0.45
2:B:145:LYS:HB3	2:B:197:THR:HB	1.98	0.45
1:E:153:SER:HB2	1:E:197:ASN:HB2	1.99	0.45
2:F:128:GLY:C	2:F:183:LYS:HB2	2.42	0.45
1:G:47:TRP:CZ2	1:G:49:GLY:HA2	2.52	0.44
2:H:76:ARG:HB3	2:H:77:ARG:HH21	1.82	0.44
2:D:184:ALA:O	2:D:188:LYS:HG2	2.18	0.44
2:H:125:LEU:O	2:H:183:LYS:HD3	2.18	0.44
1:A:199:ASN:ND2	1:A:206:LYS:HG2	2.33	0.43
2:F:186:TYR:CE2	2:F:211:ARG:HD3	2.53	0.43
1:A:4:LEU:HD22	1:A:24:THR:HG22	2.00	0.43
2:D:128:GLY:H	2:D:183:LYS:NZ	2.16	0.43
2:D:47:VAL:HA	2:D:58:VAL:HG21	2.01	0.43
1:C:169:VAL:HG11	2:D:160:GLN:HB3	2.00	0.43
2:D:35:TRP:HB2	2:D:48:ILE:HB	2.00	0.43
2:H:120:PRO:HB2	2:H:125:LEU:HD21	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:THR:HG23	1:A:210:LYS:HE3	1.99	0.42
1:A:100:SER:HB2	2:B:91:ARG:NH1	2.33	0.42
1:C:168:ALA:HA	1:C:178:LEU:HB3	2.01	0.42
2:B:183:LYS:O	2:B:187:GLU:HG2	2.19	0.42
1:G:168:ALA:HA	1:G:178:LEU:HB3	2.00	0.42
2:H:7:THR:HG23	2:H:22:SER:HB3	2.02	0.42
2:H:35:TRP:HB2	2:H:48:ILE:HB	2.02	0.42
2:F:33:LEU:HG	2:F:71:PHE:CG	2.54	0.42
2:H:46:LEU:HD21	2:H:49:TYR:HB3	2.02	0.41
1:C:11:LEU:HB2	1:C:147:PRO:HG3	2.02	0.41
2:B:40:PRO:HG3	2:B:165:GLU:HG2	2.02	0.41
1:G:82:MET:HE3	1:G:82:MET:HB2	1.84	0.41
1:C:193:THR:HG23	1:C:210:LYS:HD2	2.03	0.41
1:C:59:TYR:HB2	1:C:64:LYS:HD3	2.02	0.41
2:D:13:VAL:HG21	2:D:78:VAL:HG21	2.03	0.41
1:G:136:ALA:HB2	1:G:186:SER:HB3	2.03	0.41
2:B:120:PRO:HD3	2:B:132:VAL:HG22	2.02	0.41
1:E:126:PRO:HD3	1:E:138:LEU:HB3	2.03	0.41
2:D:186:TYR:HA	2:D:192:TYR:OH	2.20	0.41
2:H:37:LEU:HD13	2:H:86:TYR:CZ	2.56	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	216/230 (94%)	212 (98%)	4 (2%)	0	100	100
1	C	215/230 (94%)	213 (99%)	2 (1%)	0	100	100
1	E	213/230 (93%)	210 (99%)	3 (1%)	0	100	100
1	G	214/230 (93%)	211 (99%)	3 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	216/219 (99%)	213 (99%)	3 (1%)	0	100	100
2	D	217/219 (99%)	214 (99%)	3 (1%)	0	100	100
2	F	217/219 (99%)	214 (99%)	3 (1%)	0	100	100
2	H	217/219 (99%)	213 (98%)	4 (2%)	0	100	100
3	I	17/24 (71%)	17 (100%)	0	0	100	100
3	J	22/24 (92%)	22 (100%)	0	0	100	100
All	All	1764/1844 (96%)	1739 (99%)	25 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 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	183/193 (95%)	181 (99%)	2 (1%)	65	77
1	C	182/193 (94%)	181 (100%)	1 (0%)	81	88
1	E	181/193 (94%)	178 (98%)	3 (2%)	53	66
1	G	181/193 (94%)	180 (99%)	1 (1%)	78	86
2	B	195/196 (100%)	195 (100%)	0	100	100
2	D	196/196 (100%)	194 (99%)	2 (1%)	68	79
2	F	196/196 (100%)	194 (99%)	2 (1%)	68	79
2	H	196/196 (100%)	195 (100%)	1 (0%)	81	88
3	I	16/18 (89%)	16 (100%)	0	100	100
3	J	18/18 (100%)	18 (100%)	0	100	100
All	All	1544/1592 (97%)	1532 (99%)	12 (1%)	73	82

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	VAL

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Mol	Chain	Res	Type
1	A	207	VAL
1	C	207	VAL
2	D	47	VAL
2	D	127	SER
1	E	12	VAL
1	E	20	VAL
1	E	199	ASN
2	F	48	ILE
2	F	183	LYS
1	G	207	VAL
2	H	51	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	192	GLN
1	A	199	ASN
2	B	27	GLN
2	B	53	ASN
1	C	3	GLN
1	C	164	HIS
2	D	100	GLN
1	E	164	HIS
2	F	17	GLN
2	F	45	GLN
2	F	137	ASN
2	F	138	ASN
2	F	147	GLN
1	G	164	HIS
1	G	192	GLN
2	H	53	ASN
2	H	137	ASN
2	H	152	ASN
2	H	160	GLN
2	H	199	GLN
3	I	13	ASN

### 5.3.3 RNA ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	220/230 (95%)	0.10	6 (2%) 56 62	17, 25, 39, 54	0
1	C	219/230 (95%)	0.16	5 (2%) 61 66	17, 28, 39, 58	0
1	E	217/230 (94%)	0.43	6 (2%) 55 61	23, 34, 45, 59	0
1	G	218/230 (94%)	0.12	5 (2%) 61 66	15, 28, 44, 55	0
2	B	218/219 (99%)	0.12	6 (2%) 55 61	16, 28, 43, 58	0
2	D	219/219 (100%)	0.25	5 (2%) 61 66	20, 30, 49, 65	0
2	F	219/219 (100%)	0.68	18 (8%) 17 21	24, 37, 57, 73	0
2	H	219/219 (100%)	0.66	20 (9%) 15 18	19, 36, 57, 78	0
3	I	21/24 (87%)	0.93	4 (19%) 3 3	23, 33, 62, 66	0
3	J	24/24 (100%)	0.65	2 (8%) 17 20	19, 25, 57, 67	0
All	All	1794/1844 (97%)	0.33	77 (4%) 40 46	15, 30, 50, 78	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	J	10	ALA	5.5
2	F	27(C)	LEU	4.8
3	J	9	ASN	4.8
2	B	126	LYS	4.5
2	H	214	CYS	4.3
2	D	16	GLY	4.3
1	E	128	SER	4.2
2	F	214	CYS	3.8
1	C	132	SER	3.8
1	E	127	SER	3.7
1	G	128	SER	3.6
2	H	121	SER	3.5
2	F	68	GLY	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	128	SER	3.4
2	H	170	ASP	3.3
3	I	9	ASN	3.2
2	H	169	LYS	3.2
2	F	69	THR	3.2
1	E	61	ALA	3.1
2	H	168	SER	3.1
1	C	131	THR	3.1
2	H	122	ASP	3.0
1	G	133	GLY	3.0
3	I	13	ASN	3.0
2	B	16	GLY	2.9
2	B	127	SER	2.9
2	D	214	CYS	2.9
2	F	31	THR	2.9
1	C	128	SER	2.8
2	H	12	SER	2.8
2	H	80	ALA	2.8
2	F	76	ARG	2.8
3	I	24	PRO	2.7
2	H	77	ARG	2.7
2	B	129	THR	2.7
3	I	14	ALA	2.6
1	E	1	GLU	2.6
2	F	26	SER	2.6
1	A	131	THR	2.5
2	H	76	ARG	2.5
2	H	63	SER	2.5
1	G	186	SER	2.5
2	H	120	PRO	2.5
2	F	5	THR	2.4
1	A	132	SER	2.4
2	H	205	VAL	2.4
2	H	152	ASN	2.4
2	B	212	GLY	2.4
2	D	126	LYS	2.4
2	B	128	GLY	2.4
2	F	63	SER	2.4
2	F	67	SER	2.4
1	C	207	VAL	2.4
2	F	3	VAL	2.4
2	H	191	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
2	H	83	VAL	2.3
1	E	214	LYS	2.3
2	F	25	SER	2.3
1	A	101	VAL	2.3
2	H	212	GLY	2.2
2	F	184	ALA	2.2
2	F	127	SER	2.2
1	A	100	SER	2.2
1	G	41	PRO	2.2
2	F	212	GLY	2.2
1	C	116	THR	2.2
2	H	123	GLU	2.1
1	A	188	SER	2.1
1	E	134	GLY	2.1
2	F	27(D)	ASP	2.1
2	H	70	ASP	2.1
2	H	110	VAL	2.1
2	F	70	ASP	2.1
2	D	77	ARG	2.0
1	G	214	LYS	2.0
2	D	169	LYS	2.0
2	F	188	LYS	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.