



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 12, 2026 – 08:35 PM UTC

PDB ID : 6BFQ / pdb\_00006bfq  
Title : The mechanism of GM-CSF inhibition by human GM-CSF auto-antibodies  
Authors : Dhagat, U.; Hercus, T.R.; Broughton, S.E.; Nero, T.L.; Lopez, A.F.; Parker, M.W.  
Deposited on : 2017-10-26  
Resolution : 2.60 Å (reported)

This is a wwPDB X-ray Structure Validation Summary 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

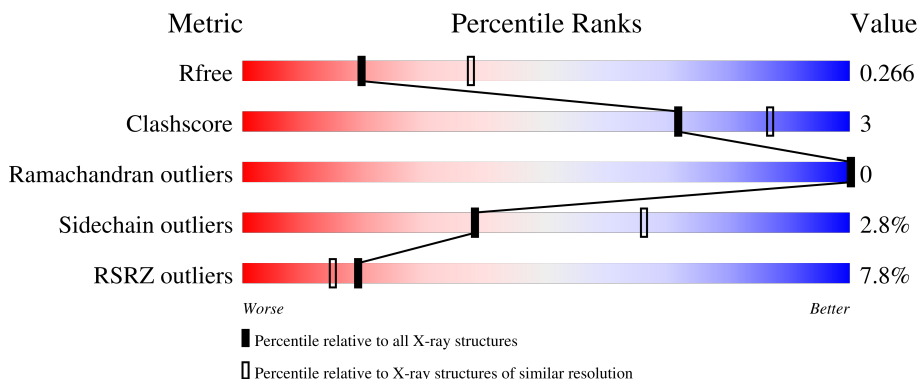
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	225	 4% 83% 9% 7%
1	C	225	 2% 88% 5% 6%
1	E	225	 2% 86% 8% 7%
1	H	225	 4% 89% 6% 5%
2	B	215	 5% 91% 7%

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Mol	Chain	Length	Quality of chain
2	D	215	<p>3% 93% 5% ..</p>
2	F	215	<p>4% 90% 8% .</p>
2	L	215	<p>2% 93% 6% .</p>
3	G	127	<p>19% 66% 6% . 26%</p>
3	I	127	<p>23% 67% 6% . 26%</p>
3	J	127	<p>24% 69% 9% . 20%</p>
3	K	127	<p>20% 70% 9% . 20%</p>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 16222 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 Fab Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	H	213	Total 1600	C 1016	N 264	O 315	S 5	0	1	0
1	A	210	Total 1571	C 999	N 257	O 310	S 5	0	0	0
1	C	211	Total 1577	C 1002	N 258	O 312	S 5	0	0	0
1	E	210	Total 1578	C 1004	N 258	O 311	S 5	0	0	0

- Molecule 2 is a protein called Fab Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	L	212	Total 1624	C 1011	N 283	O 324	S 6	0	0	0
2	B	212	Total 1624	C 1011	N 283	O 324	S 6	0	0	0
2	D	213	Total 1641	C 1021	N 287	O 327	S 6	0	1	0
2	F	213	Total 1638	C 1019	N 285	O 328	S 6	0	1	0

- Molecule 3 is a protein called Granulocyte-macrophage colony-stimulating factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	G	94	Total 762	C 487	N 125	O 142	S 8	0	0	0
3	J	102	Total 826	C 524	N 137	O 157	S 8	0	0	0
3	I	94	Total 763	C 488	N 125	O 142	S 8	0	0	0
3	K	102	Total 825	C 524	N 137	O 156	S 8	0	0	0

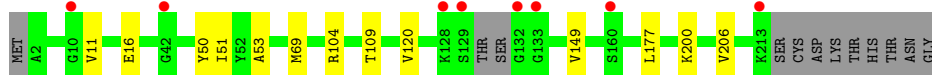
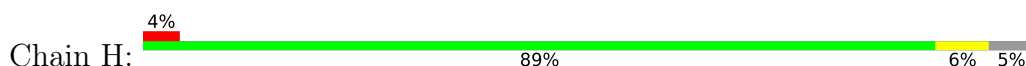
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	H	23	Total O 23 23	0	0
4	L	15	Total O 15 15	0	0
4	A	28	Total O 28 28	0	0
4	B	14	Total O 14 14	0	0
4	C	30	Total O 30 30	0	0
4	D	19	Total O 19 19	0	0
4	E	24	Total O 24 24	0	0
4	F	20	Total O 20 20	0	0
4	G	1	Total O 1 1	0	0
4	J	4	Total O 4 4	0	0
4	I	10	Total O 10 10	0	0
4	K	5	Total O 5 5	0	0

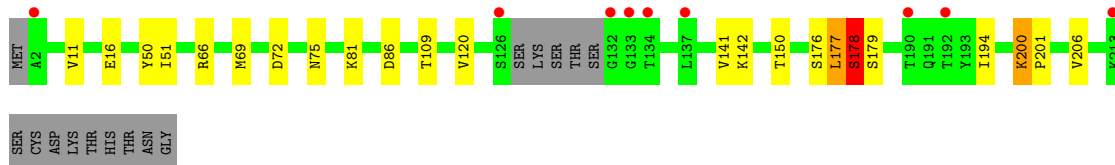
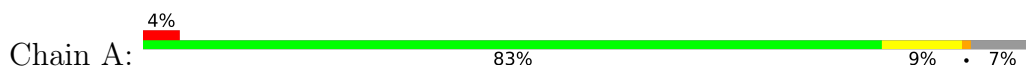
### 3 Residue-property plots [i](#)

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.

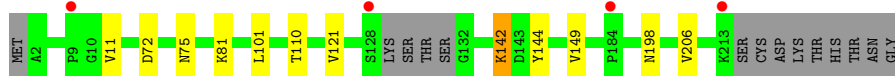
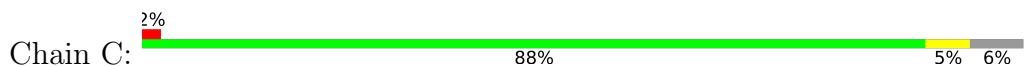
- Molecule 1: Fab Heavy chain



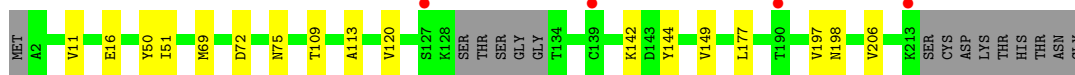
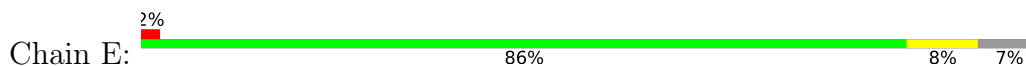
- Molecule 1: Fab Heavy chain



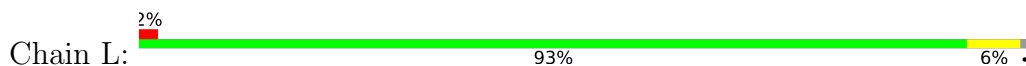
- Molecule 1: Fab Heavy chain



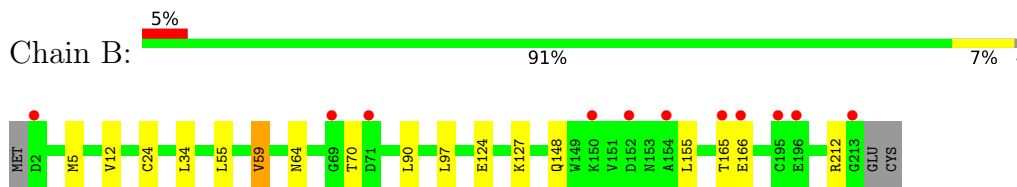
- Molecule 1: Fab Heavy chain



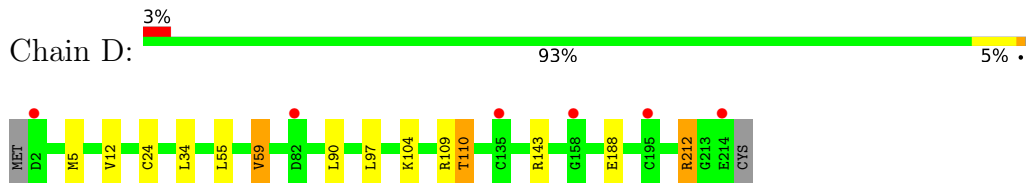
- Molecule 2: Fab Light Chain



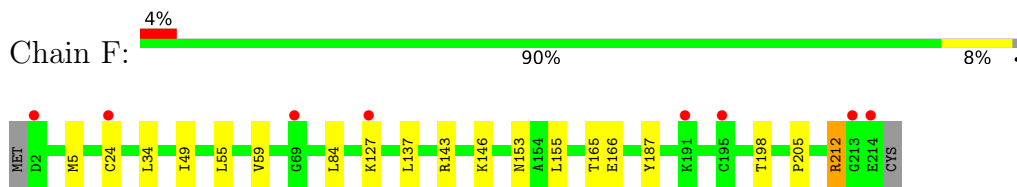
- Molecule 2: Fab Light Chain



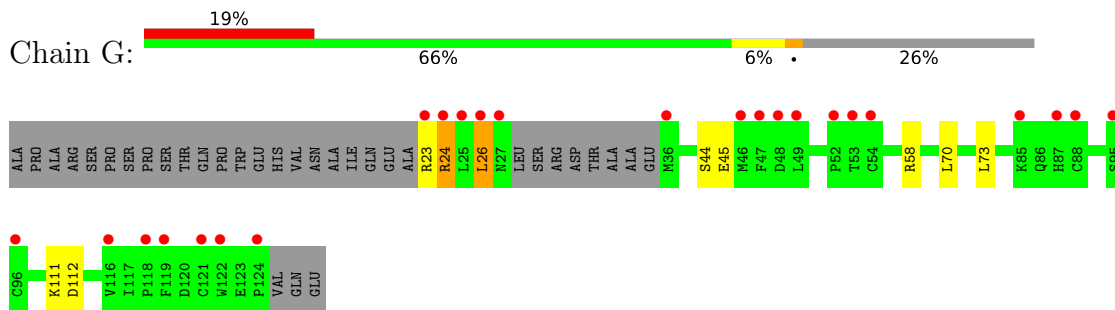
- Molecule 2: Fab Light Chain



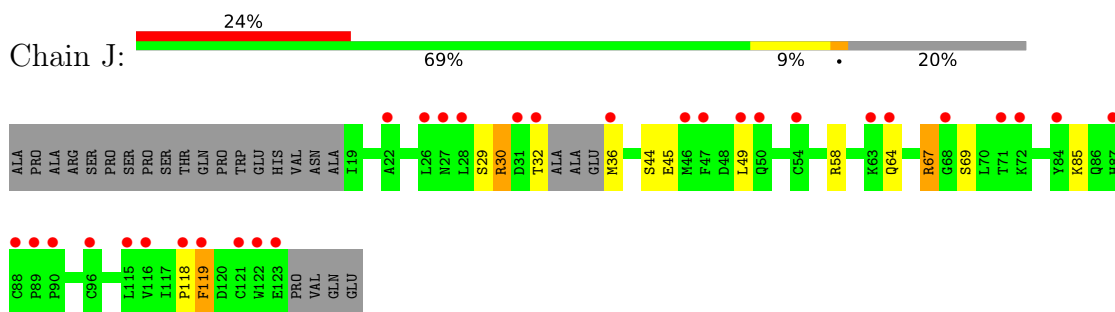
- Molecule 2: Fab Light Chain



- Molecule 3: Granulocyte-macrophage colony-stimulating factor

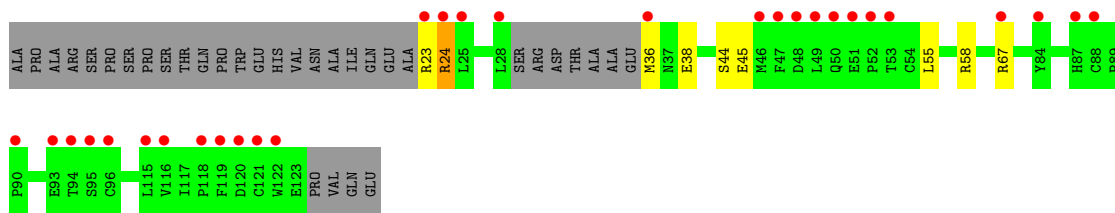


- Molecule 3: Granulocyte-macrophage colony-stimulating factor

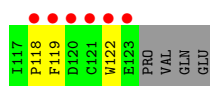
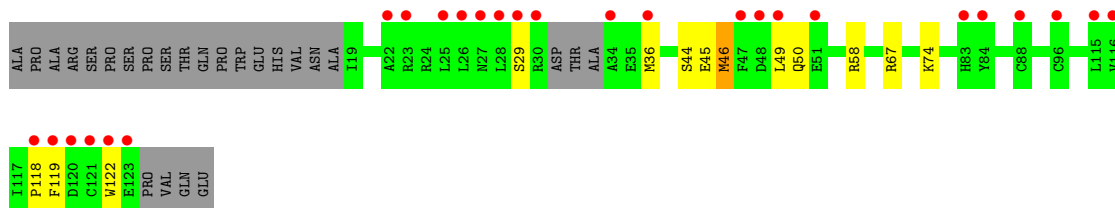


- Molecule 3: Granulocyte-macrophage colony-stimulating factor





- Molecule 3: Granulocyte-macrophage colony-stimulating factor



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	95.44Å 100.53Å 101.87Å 91.17° 117.96° 108.65°	Depositor
Resolution (Å)	50.01 – 2.60 50.01 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.6 (50.01-2.60) 98.6 (50.01-2.60)	Depositor EDS
$R_{merge}$	0.25	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.54 (at 2.61Å)	Xtrriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.230 , 0.264 0.234 , 0.266	Depositor DCC
$R_{free}$ test set	4816 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.3	Xtrriage
Anisotropy	0.438	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 33.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	16222	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.16% 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	1.00	2/1611 (0.1%)	0.98	2/2203 (0.1%)
1	C	0.99	1/1617 (0.1%)	0.99	1/2211 (0.0%)
1	E	0.95	0/1618	0.96	0/2212
1	H	0.95	0/1643	0.97	0/2244
2	B	0.85	0/1656	0.90	0/2247
2	D	0.85	0/1676	0.92	1/2273 (0.0%)
2	F	0.85	0/1673	0.91	1/2270 (0.0%)
2	L	0.86	0/1656	0.91	1/2247 (0.0%)
3	G	0.90	0/778	1.07	1/1051 (0.1%)
3	I	0.90	0/778	1.07	0/1050
3	J	0.89	0/841	1.06	0/1135
3	K	0.88	0/840	1.06	2/1133 (0.2%)
All	All	0.91	3/16387 (0.0%)	0.97	9/22276 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	101	LEU	C-N	7.30	1.43	1.33
1	A	179	SER	C-O	-5.34	1.17	1.24
1	A	178	SER	C-O	-5.01	1.17	1.23

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	K	122	TRP	N-CA-C	-5.48	101.14	108.86
1	C	142	LYS	N-CA-C	5.42	118.39	109.72
1	A	200	LYS	CA-C-N	5.32	125.39	119.32
1	A	200	LYS	C-N-CA	5.32	125.39	119.32
2	L	143	ARG	N-CA-C	5.13	116.95	111.36

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	A	1571	0	1541	15	0
1	C	1577	0	1546	4	0
1	E	1578	0	1553	12	0
1	H	1600	0	1577	11	0
2	B	1624	0	1593	8	0
2	D	1641	0	1612	6	0
2	F	1638	0	1605	12	0
2	L	1624	0	1593	7	0
3	G	762	0	761	14	0
3	I	763	0	765	5	0
3	J	826	0	824	9	0
3	K	825	0	824	6	0
4	A	28	0	0	0	0
4	B	14	0	0	0	0
4	C	30	0	0	0	0
4	D	19	0	0	0	0
4	E	24	0	0	1	0
4	F	20	0	0	0	0
4	G	1	0	0	0	0
4	H	23	0	0	0	0
4	I	10	0	0	0	0
4	J	4	0	0	2	0
4	K	5	0	0	1	0
4	L	15	0	0	0	0
All	All	16222	0	15794	105	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.

The worst 5 of 105 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)
3:G:24:ARG:HG2	3:G:24:ARG:HH21	1.39	0.83
3:G:26:LEU:HD13	3:G:70:LEU:HD22	1.65	0.76
2:F:49:ILE:HD13	2:F:55:LEU:HD12	1.68	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:LYS:HB2	1:A:201:PRO:HD3	1.69	0.75
3:G:24:ARG:HG2	3:G:24:ARG:NH2	1.98	0.74

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	206/225 (92%)	195 (95%)	11 (5%)	0	100	100
1	C	207/225 (92%)	196 (95%)	11 (5%)	0	100	100
1	E	206/225 (92%)	195 (95%)	11 (5%)	0	100	100
1	H	210/225 (93%)	200 (95%)	10 (5%)	0	100	100
2	B	210/215 (98%)	201 (96%)	9 (4%)	0	100	100
2	D	212/215 (99%)	205 (97%)	7 (3%)	0	100	100
2	F	212/215 (99%)	204 (96%)	8 (4%)	0	100	100
2	L	210/215 (98%)	202 (96%)	8 (4%)	0	100	100
3	G	90/127 (71%)	84 (93%)	6 (7%)	0	100	100
3	I	90/127 (71%)	84 (93%)	6 (7%)	0	100	100
3	J	98/127 (77%)	92 (94%)	6 (6%)	0	100	100
3	K	98/127 (77%)	92 (94%)	6 (6%)	0	100	100
All	All	2049/2268 (90%)	1950 (95%)	99 (5%)	0	100	100

There are no Ramachandran outliers to report.

### 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	A	178/192 (93%)	170 (96%)	8 (4%)	24	50
1	C	179/192 (93%)	176 (98%)	3 (2%)	53	78
1	E	180/192 (94%)	177 (98%)	3 (2%)	53	78
1	H	182/192 (95%)	179 (98%)	3 (2%)	55	79
2	B	185/188 (98%)	178 (96%)	7 (4%)	29	56
2	D	187/188 (100%)	182 (97%)	5 (3%)	39	67
2	F	187/188 (100%)	183 (98%)	4 (2%)	47	73
2	L	185/188 (98%)	184 (100%)	1 (0%)	81	92
3	G	89/116 (77%)	85 (96%)	4 (4%)	24	50
3	I	89/116 (77%)	86 (97%)	3 (3%)	32	60
3	J	96/116 (83%)	90 (94%)	6 (6%)	16	36
3	K	95/116 (82%)	90 (95%)	5 (5%)	20	43
All	All	1832/1984 (92%)	1780 (97%)	52 (3%)	38	66

5 of 52 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	142	LYS
3	G	24	ARG
3	K	50	GLN
1	E	198	ASN
2	F	127	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	191	GLN
1	E	43	GLN
3	K	64	GLN
2	D	38	GLN

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Mol	Chain	Res	Type
1	E	191	GLN

### 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	210/225 (93%)	0.18	9 (4%) 40 34	18, 34, 77, 102	0
1	C	211/225 (93%)	0.13	4 (1%) 66 61	16, 34, 75, 97	0
1	E	210/225 (93%)	0.11	4 (1%) 66 61	20, 34, 69, 95	0
1	H	213/225 (94%)	0.18	8 (3%) 44 38	17, 35, 70, 101	1 (0%)
2	B	212/215 (98%)	0.46	11 (5%) 33 27	22, 46, 72, 83	0
2	D	213/215 (99%)	0.37	6 (2%) 55 49	17, 43, 70, 92	1 (0%)
2	F	213/215 (99%)	0.30	8 (3%) 44 38	18, 42, 65, 99	1 (0%)
2	L	212/215 (98%)	0.18	4 (1%) 66 61	22, 42, 64, 77	0
3	G	94/127 (74%)	1.46	24 (25%) 1 1	27, 61, 101, 116	0
3	I	94/127 (74%)	1.32	29 (30%) 1 1	24, 64, 93, 124	0
3	J	102/127 (80%)	1.43	30 (29%) 1 1	24, 66, 92, 125	0
3	K	102/127 (80%)	1.30	26 (25%) 1 1	27, 65, 100, 114	0
All	All	2086/2268 (91%)	0.45	163 (7%) 19 15	16, 42, 82, 125	3 (0%)

The worst 5 of 163 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	G	116	VAL	7.4
1	H	132	GLY	5.8
3	G	25	LEU	5.6
3	K	22	ALA	5.5
3	I	118	PRO	5.3

### 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.