



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:53 PM GMT

PDB ID : 4Q6I  
Title : Crystal structure of murine 2D5 Fab, a potent anti-CD4 HIV-1-neutralizing antibody in complex with CD4  
Authors : Boyington, J.C.; Nabel, G.J.; Mascola, J.R.  
Deposited on : 2014-04-22  
Resolution : 3.65 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

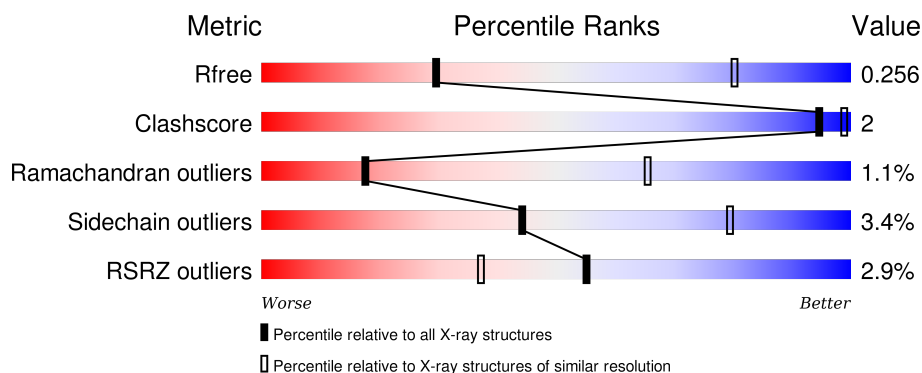
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.65 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))
$R_{free}$	91344	1010 (3.82-3.50)
Clashscore	102246	1125 (3.82-3.50)
Ramachandran outliers	100387	1079 (3.82-3.50)
Sidechain outliers	100360	1078 (3.82-3.50)
RSRZ outliers	91569	1017 (3.82-3.50)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	218	<div> <div>11%</div> <div>98%</div> <div>.</div> </div>
1	D	218	<div> <div>3%</div> <div>94%</div> <div>6%</div> </div>
1	F	218	<div> <div>3%</div> <div>94%</div> <div>6%</div> </div>
1	L	218	<div> <div>%</div> <div>94%</div> <div>6%</div> </div>
2	B	230	<div> <div>6%</div> <div>90%</div> <div>6%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
2	E	230	<div><div><div></div><div></div><div></div></div><div>2%</div><div>88%</div><div>7%</div><div></div><div></div></div>
2	G	230	<div><div><div></div><div></div><div></div></div><div>3%</div><div>87%</div><div>7%</div><div></div><div></div></div>
2	H	230	<div><div><div></div><div></div><div></div></div><div>%</div><div>87%</div><div>8%</div><div></div><div></div></div>
3	C	208	<div><div><div></div><div></div><div></div></div><div>%</div><div>79%</div><div>9%</div><div>13%</div><div></div></div>
3	I	208	<div><div><div></div><div></div><div></div></div><div></div><div>77%</div><div>11%</div><div>13%</div><div></div></div>
3	J	208	<div><div><div></div><div></div><div></div></div><div></div><div>79%</div><div>8%</div><div>13%</div><div></div></div>
3	K	208	<div><div><div></div><div></div><div></div></div><div></div><div>77%</div><div>11%</div><div>13%</div><div></div></div>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 37900 atoms, of which 18708 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 Light chain of murine 2D5 Fab.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	218	Total	C	H	N	O	S	0	0	0
			3294	1054	1598	287	346	9			
1	D	218	Total	C	H	N	O	S	0	0	0
			3293	1054	1597	287	346	9			
1	F	218	Total	C	H	N	O	S	0	0	0
			3292	1054	1596	287	346	9			
1	L	218	Total	C	H	N	O	S	0	0	0
			3294	1054	1598	287	346	9			

- Molecule 2 is a protein called Heavy chain of murine 2D5 Fab.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	222	Total	C	H	N	O	S	0	0	0
			3315	1066	1632	276	332	9			
2	E	222	Total	C	H	N	O	S	0	0	0
			3315	1066	1632	276	332	9			
2	G	222	Total	C	H	N	O	S	0	0	0
			3325	1068	1638	277	333	9			
2	H	222	Total	C	H	N	O	S	0	0	0
			3326	1068	1639	277	333	9			

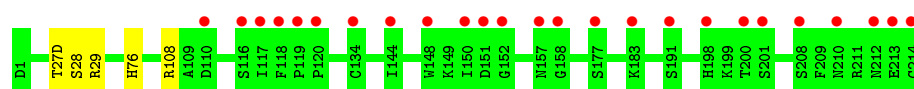
- Molecule 3 is a protein called T-cell surface glycoprotein CD4.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	182	Total	C	H	N	O	S	0	0	0
			2862	888	1445	248	277	4			
3	I	182	Total	C	H	N	O	S	0	0	0
			2862	888	1445	248	277	4			
3	J	182	Total	C	H	N	O	S	0	0	0
			2862	888	1445	248	277	4			
3	K	182	Total	C	H	N	O	S	0	0	0
			2860	888	1443	248	277	4			

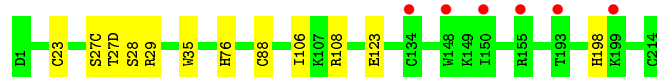
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

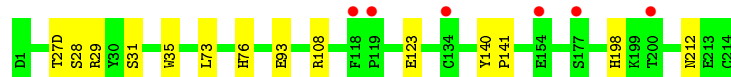
- Molecule 1: Light chain of murine 2D5 Fab



- Molecule 1: Light chain of murine 2D5 Fab



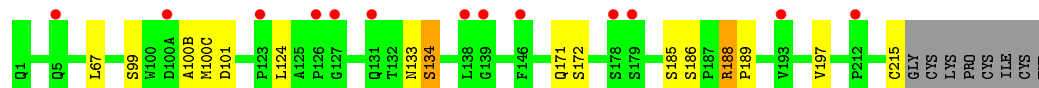
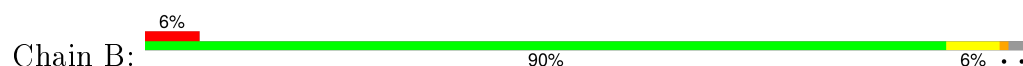
- Molecule 1: Light chain of murine 2D5 Fab



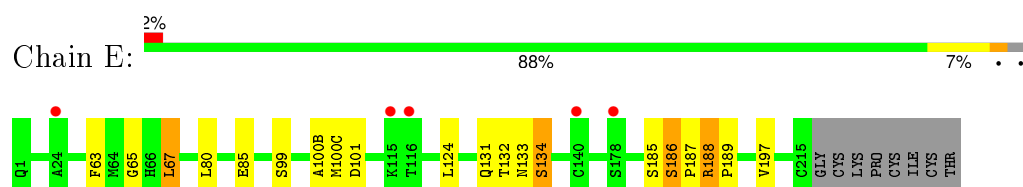
- Molecule 1: Light chain of murine 2D5 Fab



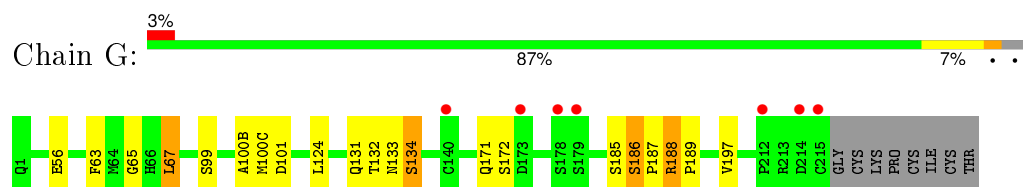
- Molecule 2: Heavy chain of murine 2D5 Fab



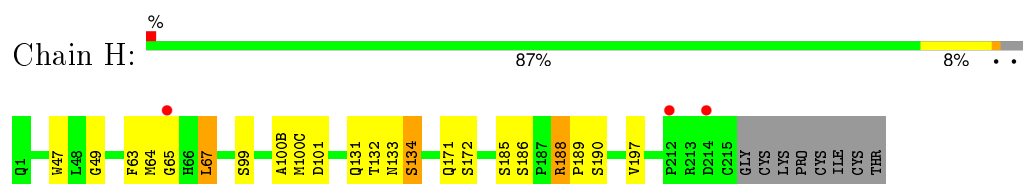
- Molecule 2: Heavy chain of murine 2D5 Fab



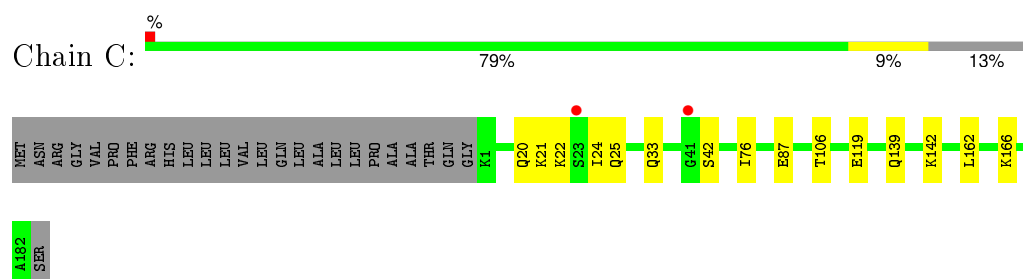
- Molecule 2: Heavy chain of murine 2D5 Fab



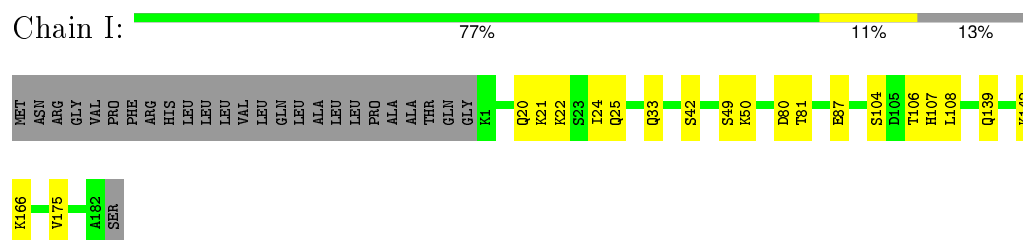
- Molecule 2: Heavy chain of murine 2D5 Fab



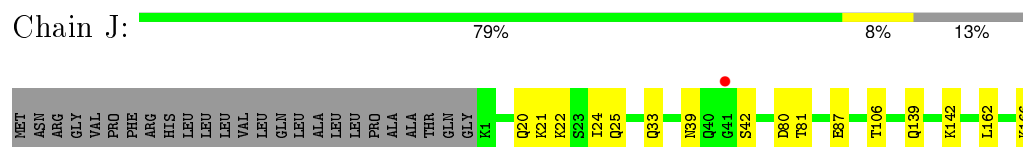
- Molecule 3: T-cell surface glycoprotein CD4



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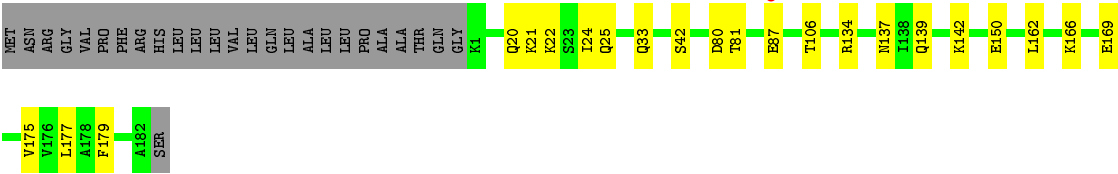
- Molecule 3: T-cell surface glycoprotein CD4

Chain K: 

77%

11%

13%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 4	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	149.55Å 149.55Å 183.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.91 – 3.65 30.02 – 3.65	Depositor EDS
% Data completeness (in resolution range)	96.8 (29.91-3.65) 96.9 (30.02-3.65)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.35 (at 3.65Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, $R_{free}$	0.226 , 0.257 0.230 , 0.256	Depositor DCC
$R_{free}$ test set	2205 reflections (5.08%)	DCC
Wilson B-factor (Å <sup>2</sup> )	129.9	Xtriage
Anisotropy	0.298	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 80.7	EDS
Estimated twinning fraction	0.024 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 43473 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	37900	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	146.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.02% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.23	0/1737	0.41	0/2359
1	D	0.23	0/1737	0.41	0/2359
1	F	0.24	0/1737	0.41	0/2359
1	L	0.23	0/1737	0.41	0/2359
2	B	0.23	0/1728	0.41	0/2359
2	E	0.23	0/1728	0.42	0/2359
2	G	0.23	0/1732	0.41	0/2364
2	H	0.24	0/1732	0.42	0/2364
3	C	0.24	0/1437	0.44	0/1937
3	I	0.24	0/1437	0.44	0/1937
3	J	0.23	0/1437	0.44	0/1937
3	K	0.24	0/1437	0.44	0/1937
All	All	0.23	0/19616	0.42	0/26630

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

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	1696	1598	1609	1	0
1	D	1696	1597	1610	5	0
1	F	1696	1596	1610	7	0
1	L	1696	1598	1610	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1683	1632	1637	4	0
2	E	1683	1632	1638	9	0
2	G	1687	1638	1644	9	0
2	H	1687	1639	1644	8	0
3	C	1417	1445	1449	4	0
3	I	1417	1445	1449	7	0
3	J	1417	1445	1449	4	0
3	K	1417	1443	1449	5	0
All	All	19192	18708	18798	64	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 2.

All (64) 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:E:134:SER:O	2:E:185:SER:N	2.21	0.74
1:A:27(D):THR:O	1:A:29:ARG:N	2.22	0.72
1:D:27(D):THR:O	1:D:29:ARG:N	2.23	0.71
1:L:27(D):THR:O	1:L:29:ARG:N	2.23	0.70
1:F:27(D):THR:O	1:F:29:ARG:N	2.24	0.70
2:B:134:SER:O	2:B:185:SER:N	2.30	0.64
2:H:134:SER:O	2:H:185:SER:N	2.31	0.63
2:G:134:SER:O	2:G:185:SER:N	2.35	0.60
2:E:100(B):ALA:O	2:E:101:ASP:N	2.36	0.59
2:B:100(B):ALA:O	2:B:101:ASP:N	2.37	0.57
2:H:100(B):ALA:O	2:H:101:ASP:N	2.38	0.57
2:G:100(B):ALA:O	2:G:101:ASP:N	2.40	0.55
1:F:212:ASN:ND2	3:K:169:GLU:O	2.34	0.54
3:J:80:ASP:OD1	3:J:81:THR:N	2.45	0.49
3:J:20:GLN:O	3:J:22:LYS:N	2.47	0.48
2:H:188:ARG:HG2	2:H:189:PRO:HA	1.96	0.48
1:F:93:GLU:OE2	3:I:50:LYS:NZ	2.47	0.48
2:G:134:SER:O	2:G:185:SER:OG	2.26	0.47
3:K:20:GLN:O	3:K:22:LYS:N	2.48	0.47
3:C:24:ILE:HG12	3:C:25:GLN:H	1.80	0.47
3:C:20:GLN:O	3:C:22:LYS:N	2.47	0.47
3:J:24:ILE:HG12	3:J:25:GLN:H	1.79	0.47
2:G:188:ARG:HG2	2:G:189:PRO:HA	1.97	0.46
3:K:24:ILE:HG12	3:K:25:GLN:H	1.80	0.46
2:E:188:ARG:HG2	2:E:189:PRO:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:131:GLN:O	2:E:132:THR:OG1	2.30	0.46
1:F:123:GLU:N	1:F:123:GLU:OE1	2.49	0.46
3:I:24:ILE:HG12	3:I:25:GLN:H	1.82	0.45
3:I:107:HIS:CE1	1:L:95:PRO:HD2	2.52	0.45
2:H:131:GLN:O	2:H:132:THR:OG1	2.29	0.45
3:C:181:LYS:NZ	2:G:56:GLU:OE1	2.50	0.44
1:L:123:GLU:N	1:L:123:GLU:OE1	2.46	0.44
1:D:123:GLU:N	1:D:123:GLU:OE1	2.47	0.44
2:H:64:MET:HA	3:I:151:LEU:HD12	2.00	0.43
2:G:131:GLN:O	2:G:132:THR:OG1	2.29	0.43
3:I:80:ASP:OD1	3:I:81:THR:N	2.50	0.43
3:J:25:GLN:NE2	3:J:39:ASN:O	2.52	0.43
3:K:134:ARG:NH1	3:K:150:GLU:OE2	2.52	0.43
2:G:186:SER:CB	2:G:187:PRO:HD3	2.49	0.42
2:E:85:GLU:N	2:E:85:GLU:OE1	2.50	0.42
2:H:171:GLN:O	2:H:172:SER:OG	2.35	0.42
2:G:65:GLY:O	2:G:67:LEU:N	2.47	0.42
2:B:171:GLN:O	2:B:172:SER:OG	2.36	0.42
3:K:80:ASP:OD1	3:K:81:THR:N	2.47	0.42
2:B:188:ARG:HG2	2:B:189:PRO:HA	2.01	0.42
2:E:65:GLY:C	2:E:67:LEU:H	2.23	0.42
1:F:198:HIS:N	1:F:198:HIS:CD2	2.88	0.42
3:C:76:ILE:HG12	3:C:119:GLU:HG2	2.02	0.41
1:L:198:HIS:CD2	1:L:198:HIS:N	2.88	0.41
2:G:171:GLN:O	2:G:172:SER:OG	2.36	0.41
2:H:47:TRP:CH2	2:H:49:GLY:HA2	2.54	0.41
1:D:198:HIS:CD2	1:D:198:HIS:N	2.88	0.41
2:H:65:GLY:O	2:H:67:LEU:N	2.45	0.41
2:E:67:LEU:HD13	2:E:80:LEU:HD11	2.03	0.41
3:I:20:GLN:O	3:I:22:LYS:N	2.54	0.41
1:D:35:TRP:CZ3	1:D:88:CYS:HB3	2.56	0.41
1:L:23:CYS:HB2	1:L:35:TRP:CH2	2.56	0.40
1:L:35:TRP:CZ3	1:L:73:LEU:HD23	2.56	0.40
1:F:35:TRP:CZ3	1:F:73:LEU:HD23	2.56	0.40
1:F:140:TYR:CG	1:F:141:PRO:HA	2.56	0.40
3:I:104:SER:HB3	3:I:108:LEU:HD21	2.04	0.40
2:E:186:SER:CB	2:E:187:PRO:HD3	2.51	0.40
2:E:134:SER:O	2:E:185:SER:OG	2.32	0.40
1:D:23:CYS:HB2	1:D:35:TRP:CH2	2.56	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	A	216/218 (99%)	202 (94%)	13 (6%)	1 (0%)	34	77
1	D	216/218 (99%)	204 (94%)	11 (5%)	1 (0%)	34	77
1	F	216/218 (99%)	203 (94%)	12 (6%)	1 (0%)	34	77
1	L	216/218 (99%)	202 (94%)	13 (6%)	1 (0%)	34	77
2	B	220/230 (96%)	196 (89%)	21 (10%)	3 (1%)	14	60
2	E	220/230 (96%)	196 (89%)	20 (9%)	4 (2%)	11	56
2	G	220/230 (96%)	196 (89%)	20 (9%)	4 (2%)	11	56
2	H	220/230 (96%)	196 (89%)	20 (9%)	4 (2%)	11	56
3	C	180/208 (86%)	159 (88%)	19 (11%)	2 (1%)	17	65
3	I	180/208 (86%)	157 (87%)	21 (12%)	2 (1%)	17	65
3	J	180/208 (86%)	159 (88%)	19 (11%)	2 (1%)	17	65
3	K	180/208 (86%)	158 (88%)	20 (11%)	2 (1%)	17	65
All	All	2464/2624 (94%)	2228 (90%)	209 (8%)	27 (1%)	17	65

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	SER
2	B	100(C)	MET
3	C	21	LYS
1	D	28	SER
2	E	100(C)	MET
1	F	28	SER
2	G	100(C)	MET
2	H	100(C)	MET
3	I	21	LYS
3	J	21	LYS
3	K	21	LYS
1	L	28	SER

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Mol	Chain	Res	Type
2	B	133	ASN
2	E	133	ASN
2	G	133	ASN
2	H	133	ASN
3	C	87	GLU
3	I	87	GLU
3	J	87	GLU
3	K	87	GLU
2	B	134	SER
2	E	134	SER
2	G	134	SER
2	E	63	PHE
2	H	134	SER
2	G	63	PHE
2	H	63	PHE

### 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	192/192 (100%)	190 (99%)	2 (1%)	82	92
1	D	192/192 (100%)	188 (98%)	4 (2%)	61	86
1	F	192/192 (100%)	189 (98%)	3 (2%)	70	89
1	L	192/192 (100%)	188 (98%)	4 (2%)	61	86
2	B	189/197 (96%)	182 (96%)	7 (4%)	41	77
2	E	189/197 (96%)	183 (97%)	6 (3%)	46	80
2	G	190/197 (96%)	184 (97%)	6 (3%)	46	80
2	H	190/197 (96%)	184 (97%)	6 (3%)	46	80
3	C	164/185 (89%)	155 (94%)	9 (6%)	27	69
3	I	164/185 (89%)	155 (94%)	9 (6%)	27	69
3	J	164/185 (89%)	156 (95%)	8 (5%)	31	71
3	K	164/185 (89%)	153 (93%)	11 (7%)	20	63

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	2182/2296 (95%)	2107 (97%)	75 (3%)	44 79

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	HIS
1	A	108	ARG
2	B	67	LEU
2	B	99	SER
2	B	124	LEU
2	B	186	SER
2	B	188	ARG
2	B	197	VAL
2	B	215	CYS
3	C	33	GLN
3	C	42	SER
3	C	106	THR
3	C	139	GLN
3	C	142	LYS
3	C	162	LEU
3	C	166	LYS
3	C	175	VAL
3	C	179	PHE
1	D	27(C)	SER
1	D	76	HIS
1	D	106	ILE
1	D	108	ARG
2	E	67	LEU
2	E	99	SER
2	E	124	LEU
2	E	186	SER
2	E	188	ARG
2	E	197	VAL
1	F	31	SER
1	F	76	HIS
1	F	108	ARG
2	G	67	LEU
2	G	99	SER
2	G	124	LEU
2	G	186	SER
2	G	188	ARG
2	G	197	VAL

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Mol	Chain	Res	Type
2	H	67	LEU
2	H	99	SER
2	H	186	SER
2	H	188	ARG
2	H	190	SER
2	H	197	VAL
3	I	33	GLN
3	I	42	SER
3	I	49	SER
3	I	106	THR
3	I	139	GLN
3	I	142	LYS
3	I	162	LEU
3	I	166	LYS
3	I	175	VAL
3	J	33	GLN
3	J	42	SER
3	J	106	THR
3	J	139	GLN
3	J	142	LYS
3	J	162	LEU
3	J	166	LYS
3	J	175	VAL
3	K	33	GLN
3	K	42	SER
3	K	106	THR
3	K	137	ASN
3	K	139	GLN
3	K	142	LYS
3	K	162	LEU
3	K	166	LYS
3	K	175	VAL
3	K	177	LEU
3	K	179	PHE
1	L	27(C)	SER
1	L	76	HIS
1	L	106	ILE
1	L	108	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	198	HIS
3	C	107	HIS
1	D	198	HIS
2	E	164	HIS
1	F	198	HIS
2	G	164	HIS
1	L	198	HIS

### 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 carbohydrates 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	218/218 (100%)	0.65	25 (11%) 6 4	124, 199, 336, 490	0
1	D	218/218 (100%)	0.22	6 (2%) 56 39	113, 179, 275, 319	0
1	F	218/218 (100%)	0.24	6 (2%) 56 39	92, 184, 297, 351	0
1	L	218/218 (100%)	0.15	3 (1%) 78 62	84, 159, 254, 372	0
2	B	222/230 (96%)	0.40	13 (5%) 26 14	107, 207, 298, 382	0
2	E	222/230 (96%)	0.08	5 (2%) 64 46	106, 166, 245, 420	0
2	G	222/230 (96%)	0.18	7 (3%) 51 34	79, 167, 244, 394	0
2	H	222/230 (96%)	0.03	3 (1%) 78 62	79, 143, 220, 391	0
3	C	182/208 (87%)	0.04	2 (1%) 82 68	101, 137, 213, 238	0
3	I	182/208 (87%)	0.11	0 100 100	93, 134, 213, 275	0
3	J	182/208 (87%)	0.06	1 (0%) 91 85	119, 163, 230, 314	0
3	K	182/208 (87%)	0.12	1 (0%) 91 85	124, 160, 224, 267	0
All	All	2488/2624 (94%)	0.20	72 (2%) 55 37	79, 163, 278, 490	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	158	GLY	6.5
1	A	152	GLY	5.9
1	A	191	SER	5.8
1	A	134	CYS	5.5
1	A	118	PHE	4.8
1	A	151	ASP	4.7
1	A	200	THR	4.6
1	A	144	ILE	4.2
1	A	214	CYS	4.0
1	A	208	SER	4.0
2	E	140	CYS	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	116	SER	3.7
1	A	110	ASP	3.6
2	B	138	LEU	3.5
1	F	134	CYS	3.5
2	B	5	GLN	3.4
1	A	119	PRO	3.4
2	H	214	ASP	3.4
1	A	157	ASN	3.3
1	L	200	THR	3.3
1	A	213	GLU	3.2
1	D	148	TRP	3.2
1	F	154	GLU	3.2
2	B	131	GLN	3.1
1	F	200	THR	3.1
2	G	140	CYS	3.1
2	B	146	PHE	3.0
1	A	148	TRP	3.0
1	D	150	ILE	3.0
2	G	178	SER	2.9
2	G	214	ASP	2.8
1	A	183	LYS	2.8
2	B	123	PRO	2.8
1	A	198	HIS	2.8
2	B	139	GLY	2.8
2	B	100(A)	ASP	2.7
3	C	23	SER	2.7
1	D	193	THR	2.7
1	A	117	ILE	2.6
2	B	179	SER	2.6
1	A	210	ASN	2.5
3	C	41	GLY	2.5
1	A	201	SER	2.5
2	E	178	SER	2.5
2	B	126	PRO	2.5
2	H	65	GLY	2.4
1	L	201	SER	2.4
2	B	212	PRO	2.4
1	D	199	LYS	2.4
1	F	119	PRO	2.4
1	D	134	CYS	2.4
3	K	87	GLU	2.4
2	B	178	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	177	SER	2.3
2	E	24	ALA	2.3
2	B	193	VAL	2.2
2	G	179	SER	2.2
1	F	118	PHE	2.2
2	G	212	PRO	2.2
1	A	150	ILE	2.1
2	B	127	GLY	2.1
3	J	41	GLY	2.1
2	E	116	THR	2.1
1	A	120	PRO	2.1
1	L	134	CYS	2.1
2	G	215	CYS	2.1
1	A	212	ASN	2.1
1	D	155	ARG	2.1
2	E	115	LYS	2.1
2	H	212	PRO	2.0
1	F	177	SER	2.0
2	G	173	ASP	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 carbohydrates 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.