



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:19 PM GMT

PDB ID : 1A5F  
Title : FAB FRAGMENT OF A MONOCLONAL ANTI-E-SELECTIN ANTIBODY  
Authors : Rodriguez-Romero, A.; Almog, O.; Tordova, M.; Randhawa, Z.  
Deposited on : 1998-02-16  
Resolution : 2.80 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

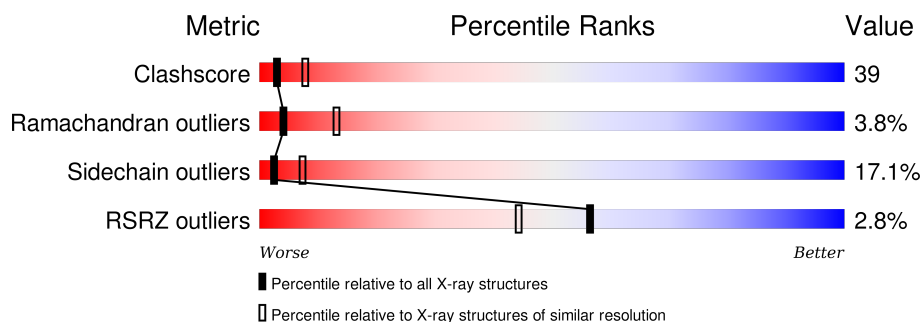
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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	L	220	 41% 50% 8%
2	H	217	 5% 34% 51% 12% ..

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3264 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 MONOCLONAL ANTI-E-SELECTIN 7A9 ANTIBODY (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	220	Total	C	N	O	S	0	0	0
			1695	1056	281	350	8			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	12	THR	SER	CONFLICT	GB 208622
L	14	THR	SER	CONFLICT	GB 208622
L	15	THR	ALA	CONFLICT	GB 208622
L	18	LYS	ARG	CONFLICT	GB 208622
L	22	THR	SER	CONFLICT	GB 208622
L	34	ALA	ASN	CONFLICT	GB 208622
L	38	TYR	PHE	CONFLICT	GB 208622
L	40	THR	ALA	CONFLICT	GB 208622
L	49	SER	PRO	CONFLICT	GB 208622
L	56	TRP	GLY	CONFLICT	GB 208622
L	80	SER	THR	CONFLICT	GB 208622
L	83	GLY	SER	CONFLICT	GB 208622
L	97	ASN	ASP	CONFLICT	GB 208622
L	98	TYR	HIS	CONFLICT	GB 208622
L	99	ASN	SER	CONFLICT	GB 208622

- Molecule 2 is a protein called MONOCLONAL ANTI-E-SELECTIN 7A9 ANTIBODY (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	212	Total	C	N	O	S	0	0	0
			1559	978	255	318	8			

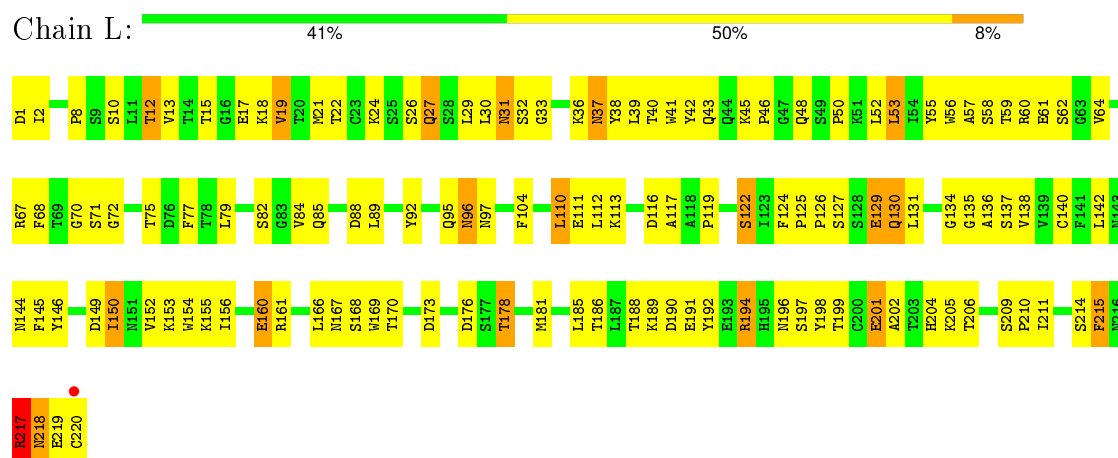
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	5	Total 5	O 5	0	0
3	L	5	Total 5	O 5	0	0

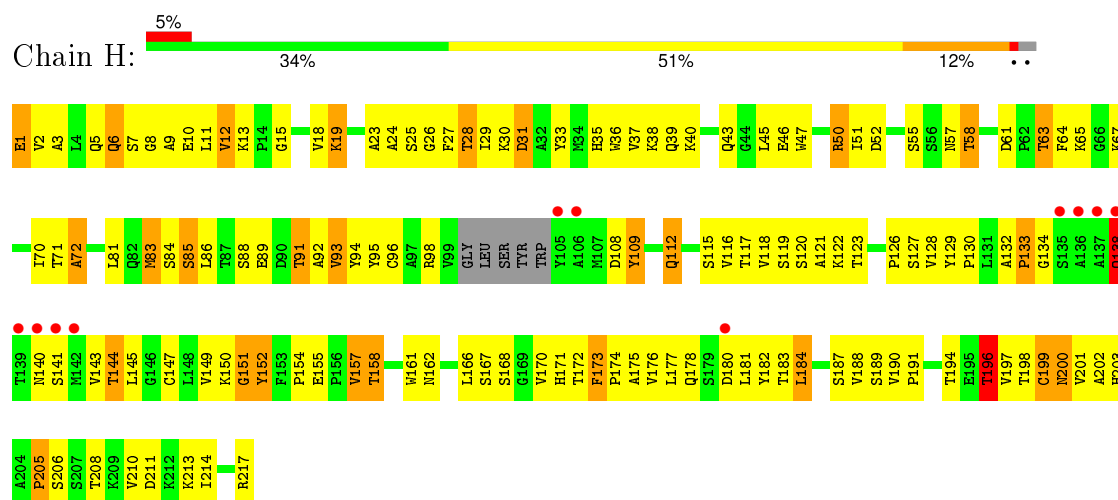
### 3 Residue-property plots

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: MONOCLONAL ANTI-E-SELECTIN 7A9 ANTIBODY (LIGHT CHAIN)



- Molecule 2: MONOCLONAL ANTI-E-SELECTIN 7A9 ANTIBODY (HEAVY CHAIN)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	44.46 Å 83.79 Å 132.45 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.00 – 2.80 35.41 – 2.81	Depositor EDS
% Data completeness (in resolution range)	86.6 (7.00-2.80) 79.0 (35.41-2.81)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.83 (at 2.81 Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, $R_{free}$	0.195 , (Not available) 0.218 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	25.2	Xtriage
Anisotropy	0.980	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 100.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 10007 reflections	Xtriage
$F_o, F_c$ correlation	0.86	EDS
Total number of atoms	3264	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.64% 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	L	0.69	0/1732	0.92	0/2354
2	H	0.73	0/1594	1.00	3/2178 (0.1%)
All	All	0.71	0/3326	0.96	3/4532 (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	L	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
2	H	138	GLN	N-CA-C	-6.30	93.99	111.00
2	H	196	THR	N-CA-C	5.83	126.75	111.00
2	H	98	ARG	N-CA-C	-5.66	95.72	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	L	38	TYR	Sidechain

### 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	L	1695	0	1622	128	0
2	H	1559	0	1507	132	0
3	H	5	0	0	0	0
3	L	5	0	0	0	0
All	All	3264	0	3129	248	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 39.

All (248) 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:L:153:LYS:HB3	1:L:201:GLU:HB3	1.43	1.01
2:H:123:THR:HG22	2:H:154:PRO:HD3	1.41	1.00
2:H:128:VAL:HG21	2:H:210:VAL:HG21	1.43	0.99
2:H:175:ALA:HB2	2:H:184:LEU:HD12	1.44	0.98
2:H:83:MET:HB3	2:H:86:LEU:HD21	1.48	0.94
1:L:40:THR:HB	1:L:95:GLN:HG2	1.52	0.92
2:H:39:GLN:HB3	2:H:93:VAL:HG23	1.53	0.88
2:H:201:VAL:HB	2:H:210:VAL:HG23	1.57	0.83
2:H:166:LEU:HD23	2:H:188:VAL:HG21	1.63	0.79
2:H:170:VAL:HG22	2:H:188:VAL:HG23	1.62	0.79
2:H:2:VAL:HG11	2:H:109:TYR:CE1	2.18	0.79
2:H:133:PRO:HG2	2:H:134:GLY:H	1.48	0.78
1:L:131:LEU:HD21	1:L:136:ALA:HB2	1.64	0.78
2:H:12:VAL:HG21	2:H:86:LEU:HD12	1.64	0.77
1:L:166:LEU:HD13	2:H:176:VAL:HG11	1.66	0.77
1:L:131:LEU:HD22	1:L:189:LYS:HG3	1.64	0.77
1:L:137:SER:HA	1:L:186:THR:HA	1.67	0.76
2:H:152:TYR:CE1	2:H:182:TYR:HB2	2.22	0.75
2:H:2:VAL:HG11	2:H:109:TYR:CD1	2.22	0.75
1:L:60:ARG:HG2	1:L:64:VAL:HB	1.69	0.75
2:H:40:LYS:HE2	2:H:43:GLN:OE1	1.88	0.74
1:L:2:ILE:HG12	1:L:27:GLN:HB2	1.70	0.72
1:L:2:ILE:HB	1:L:96:ASN:HD22	1.52	0.72
2:H:152:TYR:HE2	2:H:157:VAL:HB	1.55	0.71
1:L:166:LEU:HD13	2:H:176:VAL:CG1	2.20	0.71
2:H:15:GLY:O	2:H:85:SER:HA	1.92	0.70
1:L:40:THR:HG23	1:L:52:LEU:HD13	1.73	0.70
1:L:131:LEU:CD2	1:L:136:ALA:HB2	2.23	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:161:ARG:CG	1:L:185:LEU:HD11	2.22	0.69
1:L:142:LEU:HB2	1:L:181:MET:HG2	1.74	0.69
2:H:121:ALA:HB2	2:H:180:ASP:HB3	1.74	0.68
2:H:108:ASP:O	2:H:109:TYR:CG	2.48	0.67
2:H:196:THR:HG22	2:H:213:LYS:HE3	1.76	0.67
1:L:161:ARG:HG3	1:L:185:LEU:HD11	1.76	0.67
2:H:177:LEU:HD12	2:H:181:LEU:O	1.94	0.66
1:L:125:PRO:HD2	2:H:217:ARG:HH21	1.60	0.66
2:H:24:ALA:HB1	2:H:27:PHE:CE1	2.30	0.66
2:H:6:GLN:HB2	2:H:112:GLN:HG3	1.77	0.66
1:L:50:PRO:HG2	2:H:45:LEU:HD11	1.78	0.66
1:L:13:VAL:CG2	1:L:84:VAL:HG11	2.26	0.65
1:L:127:SER:OG	1:L:129:GLU:HG2	1.97	0.64
1:L:24:LYS:HA	1:L:75:THR:O	1.98	0.64
2:H:24:ALA:HB1	2:H:27:PHE:CZ	2.33	0.64
2:H:171:HIS:HB2	2:H:187:SER:OG	1.98	0.63
1:L:116:ASP:OD2	1:L:205:LYS:HD3	1.99	0.62
1:L:156:ILE:HG23	1:L:198:TYR:CE1	2.34	0.62
2:H:51:ILE:HB	2:H:70:ILE:HG22	1.80	0.62
2:H:175:ALA:HB1	2:H:182:TYR:HB3	1.80	0.61
2:H:198:THR:HG22	2:H:199:CYS:N	2.15	0.61
1:L:126:PRO:HG3	1:L:137:SER:O	2.01	0.61
1:L:126:PRO:HD3	1:L:138:VAL:HG22	1.81	0.61
2:H:151:GLY:HA2	2:H:181:LEU:HB3	1.82	0.60
1:L:2:ILE:HA	1:L:26:SER:HG	1.67	0.60
2:H:88:SER:O	2:H:91:THR:HB	2.02	0.60
2:H:47:TRP:HZ2	2:H:50:ARG:HG2	1.67	0.60
1:L:45:LYS:HB2	1:L:48:GLN:HG3	1.84	0.59
1:L:88:ASP:O	1:L:110:LEU:HD23	2.03	0.59
1:L:153:LYS:HZ2	1:L:160:GLU:HG3	1.68	0.59
1:L:40:THR:HG21	1:L:42:TYR:CE2	2.38	0.59
2:H:152:TYR:CE2	2:H:157:VAL:HB	2.38	0.58
2:H:130:PRO:CB	2:H:214:ILE:HD12	2.33	0.58
1:L:129:GLU:HG3	2:H:129:TYR:HD1	1.68	0.58
1:L:72:GLY:HA3	1:L:77:PHE:HA	1.84	0.58
1:L:204:HIS:ND1	1:L:205:LYS:N	2.51	0.58
2:H:33:TYR:CG	2:H:33:TYR:O	2.57	0.58
2:H:130:PRO:HB3	2:H:214:ILE:HD12	1.86	0.57
2:H:122:LYS:HE2	2:H:123:THR:OG1	2.04	0.57
2:H:175:ALA:HB2	2:H:184:LEU:CD1	2.27	0.57
2:H:40:LYS:HA	2:H:92:ALA:HB1	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:144:ASN:HD22	1:L:178:THR:HG21	1.70	0.57
2:H:38:LYS:NZ	2:H:94:TYR:OH	2.38	0.56
2:H:39:GLN:O	2:H:92:ALA:HB1	2.05	0.56
1:L:2:ILE:HB	1:L:96:ASN:ND2	2.20	0.56
2:H:61:ASP:OD1	2:H:63:THR:HG23	2.05	0.56
1:L:37:ASN:H	1:L:37:ASN:HD22	1.52	0.56
1:L:190:ASP:O	1:L:194:ARG:HB2	2.06	0.56
2:H:50:ARG:NH2	2:H:52:ASP:HB2	2.20	0.56
1:L:119:PRO:HG2	1:L:211:ILE:CD1	2.36	0.56
1:L:124:PHE:HB3	2:H:217:ARG:NH2	2.19	0.56
1:L:52:LEU:HB3	2:H:108:ASP:HB2	1.88	0.55
2:H:33:TYR:HE2	2:H:35:HIS:CE1	2.24	0.55
2:H:51:ILE:HG13	2:H:58:THR:HB	1.89	0.55
2:H:150:LYS:HG3	2:H:183:THR:OG1	2.07	0.55
2:H:128:VAL:HG21	2:H:210:VAL:CG2	2.27	0.55
2:H:152:TYR:CD1	2:H:152:TYR:N	2.75	0.55
2:H:176:VAL:O	2:H:182:TYR:HA	2.07	0.55
1:L:89:LEU:HD11	1:L:112:LEU:HG	1.88	0.55
1:L:144:ASN:HA	1:L:178:THR:HG23	1.88	0.55
1:L:131:LEU:O	1:L:189:LYS:HD2	2.06	0.55
1:L:125:PRO:HB3	1:L:215:PHE:CE1	2.41	0.54
2:H:11:LEU:HD12	2:H:117:THR:HB	1.88	0.54
2:H:12:VAL:HG22	2:H:13:LYS:O	2.06	0.54
1:L:116:ASP:HB3	1:L:206:THR:HG22	1.88	0.54
1:L:18:LYS:HA	1:L:82:SER:HA	1.89	0.54
1:L:196:ASN:O	1:L:217:ARG:HB2	2.07	0.54
1:L:153:LYS:NZ	1:L:160:GLU:HG3	2.22	0.54
1:L:58:SER:HA	1:L:70:GLY:HA3	1.89	0.53
1:L:167:ASN:HB3	1:L:181:MET:CE	2.39	0.53
1:L:40:THR:HG22	1:L:41:TRP:N	2.22	0.53
2:H:161:TRP:HE3	2:H:198:THR:O	1.92	0.53
1:L:17:GLU:O	1:L:84:VAL:HG12	2.10	0.52
1:L:40:THR:HG23	1:L:52:LEU:CD1	2.40	0.52
2:H:152:TYR:N	2:H:152:TYR:HD1	2.08	0.52
2:H:167:SER:O	2:H:170:VAL:HB	2.10	0.52
1:L:12:THR:HG22	1:L:113:LYS:HG3	1.91	0.52
1:L:42:TYR:CD1	1:L:50:PRO:HB2	2.45	0.52
1:L:117:ALA:HB3	1:L:145:PHE:HA	1.91	0.51
2:H:190:VAL:HG23	2:H:191:PRO:O	2.11	0.51
1:L:130:GLN:HA	1:L:130:GLN:HE21	1.75	0.51
1:L:125:PRO:HG2	2:H:217:ARG:NE	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:39:LEU:HD22	1:L:77:PHE:CG	2.46	0.51
1:L:13:VAL:HG21	1:L:84:VAL:HG11	1.92	0.51
1:L:119:PRO:HG2	1:L:211:ILE:HD12	1.93	0.51
1:L:31:ASN:O	1:L:33:GLY:N	2.44	0.51
2:H:38:LYS:HD3	2:H:46:GLU:OE2	2.11	0.51
1:L:142:LEU:HD23	1:L:150:ILE:HD11	1.92	0.51
1:L:40:THR:CG2	1:L:41:TRP:N	2.74	0.50
2:H:162:ASN:HD22	2:H:166:LEU:HD22	1.77	0.50
2:H:3:ALA:HB3	2:H:25:SER:OG	2.10	0.50
1:L:13:VAL:HG23	1:L:19:VAL:HG23	1.93	0.50
2:H:6:GLN:HB2	2:H:112:GLN:OE1	2.11	0.50
1:L:142:LEU:HD12	1:L:142:LEU:N	2.27	0.50
1:L:116:ASP:HA	1:L:146:TYR:O	2.12	0.49
1:L:37:ASN:HB2	1:L:57:ALA:HB2	1.94	0.49
2:H:126:PRO:CA	2:H:152:TYR:HB3	2.42	0.49
1:L:95:GLN:HA	1:L:104:PHE:HA	1.93	0.49
2:H:133:PRO:HG2	2:H:134:GLY:N	2.22	0.49
1:L:173:ASP:HB3	1:L:178:THR:H	1.78	0.49
1:L:29:LEU:HD12	1:L:39:LEU:HB2	1.93	0.49
2:H:151:GLY:N	2:H:182:TYR:O	2.45	0.49
1:L:199:THR:HA	1:L:214:SER:HA	1.94	0.49
1:L:168:SER:OG	2:H:174:PRO:HD2	2.12	0.49
2:H:128:VAL:CG2	2:H:210:VAL:HG21	2.31	0.49
1:L:111:GLU:O	1:L:111:GLU:HG3	2.12	0.49
2:H:198:THR:CG2	2:H:199:CYS:N	2.76	0.49
1:L:97:ASN:O	1:L:97:ASN:ND2	2.45	0.48
2:H:172:THR:HG23	2:H:172:THR:O	2.13	0.48
1:L:96:ASN:C	1:L:96:ASN:OD1	2.52	0.48
2:H:5:GLN:HB2	2:H:23:ALA:HB3	1.94	0.48
2:H:47:TRP:HZ2	2:H:50:ARG:CG	2.25	0.48
1:L:219:GLU:HA	1:L:219:GLU:OE1	2.13	0.48
1:L:124:PHE:HB3	2:H:217:ARG:HH22	1.77	0.48
2:H:50:ARG:NH2	2:H:52:ASP:CB	2.77	0.48
2:H:12:VAL:HG13	2:H:118:VAL:HG22	1.95	0.48
1:L:52:LEU:HG	1:L:61:GLU:HG3	1.94	0.48
2:H:29:ILE:HG13	2:H:30:LYS:N	2.27	0.48
1:L:125:PRO:HD2	2:H:217:ARG:NH2	2.28	0.48
1:L:126:PRO:HG2	1:L:136:ALA:HB1	1.96	0.47
1:L:53:LEU:HA	1:L:64:VAL:HG21	1.96	0.47
1:L:167:ASN:HB2	1:L:169:TRP:CH2	2.50	0.47
2:H:8:GLY:N	2:H:112:GLN:OE1	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:51:ILE:HA	2:H:58:THR:HB	1.96	0.47
1:L:95:GLN:HB3	1:L:104:PHE:CD1	2.49	0.47
2:H:71:THR:O	2:H:72:ALA:HB2	2.14	0.47
2:H:188:VAL:O	2:H:188:VAL:HG13	2.14	0.47
1:L:150:ILE:HD12	1:L:204:HIS:HB2	1.97	0.47
2:H:132:ALA:O	2:H:217:ARG:CZ	2.63	0.47
2:H:19:LYS:O	2:H:19:LYS:CE	2.62	0.47
2:H:9:ALA:HA	2:H:115:SER:O	2.15	0.47
1:L:122:SER:O	1:L:140:CYS:HA	2.15	0.47
1:L:40:THR:HB	1:L:95:GLN:CG	2.34	0.47
2:H:201:VAL:HB	2:H:210:VAL:CG2	2.38	0.46
1:L:153:LYS:NZ	1:L:160:GLU:CG	2.78	0.46
2:H:11:LEU:HD11	2:H:119:SER:HB3	1.96	0.46
2:H:61:ASP:HB3	2:H:64:PHE:HD2	1.81	0.46
1:L:209:SER:HB2	1:L:210:PRO:HD2	1.97	0.46
2:H:51:ILE:HG23	2:H:51:ILE:O	2.14	0.46
2:H:19:LYS:HA	2:H:81:LEU:O	2.16	0.46
1:L:21:MET:SD	1:L:92:TYR:HB2	2.55	0.46
1:L:55:TYR:CE2	1:L:61:GLU:HG2	2.51	0.46
2:H:47:TRP:CZ2	2:H:50:ARG:HG2	2.49	0.46
1:L:145:PHE:CE2	1:L:150:ILE:HG21	2.51	0.45
2:H:52:ASP:N	2:H:57:ASN:O	2.47	0.45
2:H:19:LYS:O	2:H:19:LYS:HE2	2.16	0.45
2:H:12:VAL:HG12	2:H:116:VAL:HG22	1.98	0.45
1:L:43:GLN:HB2	1:L:53:LEU:HD21	1.99	0.45
1:L:145:PHE:CE2	1:L:150:ILE:CG2	2.99	0.45
2:H:132:ALA:O	2:H:217:ARG:NH2	2.49	0.45
1:L:149:ASP:C	1:L:150:ILE:HG22	2.38	0.45
1:L:194:ARG:HD3	1:L:194:ARG:O	2.17	0.45
2:H:28:THR:HB	2:H:31:ASP:OD1	2.16	0.45
1:L:42:TYR:CD1	1:L:50:PRO:CB	3.00	0.44
2:H:40:LYS:HA	2:H:92:ALA:CB	2.47	0.44
1:L:2:ILE:HA	1:L:26:SER:OG	2.17	0.44
1:L:176:ASP:OD2	1:L:178:THR:HB	2.16	0.44
1:L:37:ASN:H	1:L:37:ASN:ND2	2.14	0.44
2:H:173:PHE:HA	2:H:174:PRO:HD3	1.81	0.44
1:L:72:GLY:CA	1:L:77:PHE:HA	2.47	0.44
2:H:143:VAL:HB	2:H:190:VAL:O	2.17	0.44
2:H:35:HIS:HD2	2:H:47:TRP:HE1	1.65	0.44
2:H:170:VAL:HG12	2:H:171:HIS:N	2.33	0.44
2:H:33:TYR:CD2	2:H:33:TYR:O	2.71	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:89:LEU:HD11	1:L:112:LEU:CG	2.48	0.44
1:L:218:ASN:HA	1:L:218:ASN:HD22	1.54	0.44
2:H:197:VAL:O	2:H:214:ILE:HG12	2.18	0.43
2:H:138:GLN:C	2:H:140:ASN:H	2.20	0.43
1:L:142:LEU:HD21	1:L:202:ALA:HB2	2.00	0.43
2:H:126:PRO:HB3	2:H:152:TYR:HB3	2.00	0.43
2:H:157:VAL:HG23	2:H:203:HIS:CD2	2.53	0.43
1:L:68:PHE:HD1	1:L:79:LEU:HD11	1.82	0.43
2:H:11:LEU:HA	2:H:117:THR:O	2.18	0.43
2:H:1:GLU:C	2:H:26:GLY:HA3	2.39	0.43
2:H:177:LEU:HD13	2:H:182:TYR:CE2	2.54	0.42
2:H:108:ASP:O	2:H:109:TYR:CD1	2.72	0.42
1:L:40:THR:CG2	1:L:52:LEU:HD13	2.45	0.42
2:H:88:SER:O	2:H:91:THR:N	2.45	0.42
1:L:67:ARG:CZ	1:L:85:GLN:HG3	2.49	0.42
1:L:154:TRP:HB3	1:L:185:LEU:CD1	2.49	0.42
1:L:192:TYR:O	1:L:198:TYR:OH	2.36	0.42
2:H:11:LEU:HD12	2:H:117:THR:O	2.19	0.42
1:L:154:TRP:O	1:L:160:GLU:HG2	2.20	0.42
2:H:203:HIS:HB3	2:H:208:THR:OG1	2.20	0.42
2:H:1:GLU:HG2	2:H:2:VAL:N	2.33	0.42
2:H:38:LYS:HE3	2:H:40:LYS:HD2	2.02	0.42
1:L:130:GLN:O	1:L:135:GLY:O	2.37	0.42
2:H:161:TRP:CE3	2:H:198:THR:O	2.72	0.42
2:H:18:VAL:HG12	2:H:86:LEU:HD11	2.01	0.42
1:L:167:ASN:HB2	1:L:169:TRP:CZ3	2.54	0.42
1:L:144:ASN:ND2	1:L:178:THR:HG21	2.32	0.42
1:L:56:TRP:O	1:L:57:ALA:HB3	2.20	0.42
1:L:155:LYS:HB3	1:L:160:GLU:HA	2.02	0.42
2:H:38:LYS:HB3	2:H:46:GLU:HG3	2.02	0.42
1:L:196:ASN:HA	1:L:217:ARG:HB2	2.02	0.42
2:H:200:ASN:HB3	2:H:211:ASP:OD1	2.19	0.42
2:H:154:PRO:HD2	2:H:205:PRO:CG	2.49	0.42
2:H:157:VAL:HG12	2:H:184:LEU:HD13	2.02	0.42
1:L:45:LYS:HB2	1:L:48:GLN:CG	2.49	0.41
1:L:144:ASN:HA	1:L:178:THR:CG2	2.50	0.41
1:L:8:PRO:HB2	1:L:10:SER:O	2.19	0.41
2:H:7:SER:N	2:H:112:GLN:OE1	2.53	0.41
2:H:67:LYS:HD3	2:H:84:SER:O	2.20	0.41
2:H:144:THR:HA	2:H:189:SER:HA	2.02	0.41
1:L:45:LYS:O	1:L:48:GLN:HG2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:12:THR:HG23	1:L:111:GLU:OE2	2.20	0.41
1:L:53:LEU:O	1:L:64:VAL:HG21	2.20	0.41
2:H:149:VAL:O	2:H:183:THR:HA	2.19	0.41
1:L:217:ARG:NH1	1:L:217:ARG:HB3	2.35	0.41
2:H:83:MET:HB3	2:H:86:LEU:CD2	2.34	0.41
2:H:158:THR:HG22	2:H:202:ALA:HB3	2.03	0.41
1:L:119:PRO:HG3	1:L:150:ILE:HD12	2.02	0.40
1:L:125:PRO:HG2	2:H:217:ARG:CZ	2.51	0.40
1:L:125:PRO:HB3	1:L:215:PHE:CZ	2.55	0.40
2:H:133:PRO:CG	2:H:134:GLY:N	2.83	0.40
1:L:198:TYR:CD2	1:L:215:PHE:CZ	3.10	0.40
2:H:37:VAL:HB	2:H:95:TYR:HB2	2.03	0.40
1:L:55:TYR:CE1	1:L:59:THR:HG21	2.57	0.40
2:H:36:TRP:CE3	2:H:81:LEU:HD22	2.57	0.40
1:L:36:LYS:HB3	1:L:36:LYS:HE2	1.80	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	L	218/220 (99%)	189 (87%)	23 (11%)	6 (3%)	6	21
2	H	208/217 (96%)	159 (76%)	39 (19%)	10 (5%)	3	9
All	All	426/437 (98%)	348 (82%)	62 (15%)	16 (4%)	4	13

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	62	SER
2	H	109	TYR
2	H	133	PRO

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Mol	Chain	Res	Type
2	H	196	THR
1	L	32	SER
1	L	134	GLY
2	H	65	LYS
2	H	151	GLY
2	H	194	THR
1	L	46	PRO
1	L	217	ARG
2	H	89	GLU
2	H	205	PRO
1	L	150	ILE
2	H	72	ALA
2	H	85	SER

### 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	L	194/196 (99%)	165 (85%)	29 (15%)	4	11
2	H	174/183 (95%)	140 (80%)	34 (20%)	2	5
All	All	368/379 (97%)	305 (83%)	63 (17%)	2	7

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	1	ASP
1	L	12	THR
1	L	15	THR
1	L	19	VAL
1	L	22	THR
1	L	27	GLN
1	L	30	LEU
1	L	31	ASN
1	L	37	ASN
1	L	53	LEU

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Mol	Chain	Res	Type
1	L	71	SER
1	L	96	ASN
1	L	110	LEU
1	L	122	SER
1	L	129	GLU
1	L	130	GLN
1	L	152	VAL
1	L	160	GLU
1	L	170	THR
1	L	178	THR
1	L	188	THR
1	L	191	GLU
1	L	194	ARG
1	L	197	SER
1	L	201	GLU
1	L	215	PHE
1	L	217	ARG
1	L	218	ASN
1	L	220	CYS
2	H	1	GLU
2	H	6	GLN
2	H	10	GLU
2	H	12	VAL
2	H	19	LYS
2	H	28	THR
2	H	31	ASP
2	H	50	ARG
2	H	55	SER
2	H	58	THR
2	H	63	THR
2	H	83	MET
2	H	91	THR
2	H	93	VAL
2	H	96	CYS
2	H	112	GLN
2	H	120	SER
2	H	127	SER
2	H	138	GLN
2	H	141	SER
2	H	144	THR
2	H	145	LEU
2	H	147	CYS

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Mol	Chain	Res	Type
2	H	152	TYR
2	H	155	GLU
2	H	157	VAL
2	H	158	THR
2	H	168	SER
2	H	173	PHE
2	H	178	GLN
2	H	184	LEU
2	H	199	CYS
2	H	200	ASN
2	H	206	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	31	ASN
1	L	37	ASN
1	L	85	GLN
1	L	144	ASN
1	L	218	ASN
2	H	82	GLN
2	H	162	ASN
2	H	171	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

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, 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	L	220/220 (100%)	-0.28	1 (0%) 91 88	2, 11, 30, 59	0
2	H	212/217 (97%)	-0.04	11 (5%) 31 20	2, 12, 43, 62	0
All	All	432/437 (98%)	-0.16	12 (2%) 56 44	2, 12, 39, 62	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	140	ASN	7.9
2	H	139	THR	5.9
2	H	141	SER	4.5
2	H	137	ALA	4.2
2	H	142	MET	3.5
2	H	136	ALA	3.1
1	L	220	CYS	2.7
2	H	180	ASP	2.6
2	H	138	GLN	2.4
2	H	106	ALA	2.2
2	H	105	TYR	2.1
2	H	135	SER	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates

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.