



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:48 PM GMT

PDB ID : 1CL7  
Title : ANTI HIV1 PROTEASE FAB  
Authors : Lescar, J.; Bentley, G.A.  
Deposited on : 1999-05-06  
Resolution : 3.00 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) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

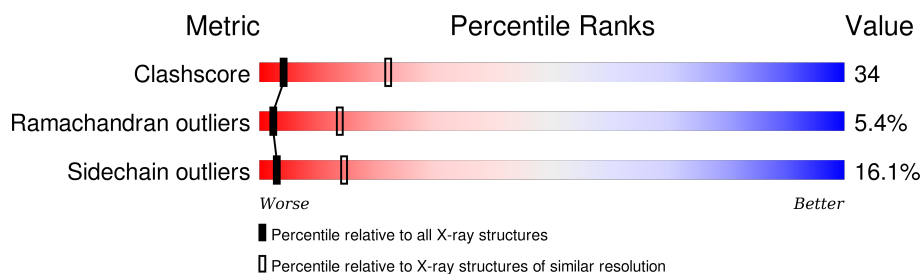
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	216	<div> <div></div> <div>39%</div> <div>51%</div> <div>10%</div> </div>
2	H	135	<div> <div></div> <div>47%</div> <div>37%</div> <div>16%</div> <div>.</div> </div>
3	I	82	<div> <div></div> <div>27%</div> <div>55%</div> <div>18%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3355 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 PROTEIN (IGG1 ANTIBODY 1696 (light chain)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	216	Total	C	N	O	S	0	0	0
			1686	1058	284	338	6			

- Molecule 2 is a protein called PROTEIN (IGG1 ANTIBODY 1696 (variable heavy chain)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	135	Total	C	N	O	S	0	0	0
			1063	676	172	212	3			

- Molecule 3 is a protein called PROTEIN (IGG1 ANTIBODY 1696 (constant heavy chain)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	82	Total	C	N	O	S	0	0	0
			606	381	100	121	4			

### 3 Residue-property plots

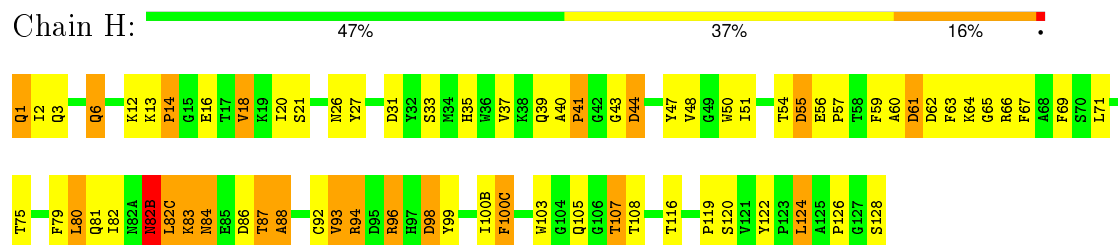
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ( $\text{RSRZ} > 2$ ). 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.

Note EDS was not executed.

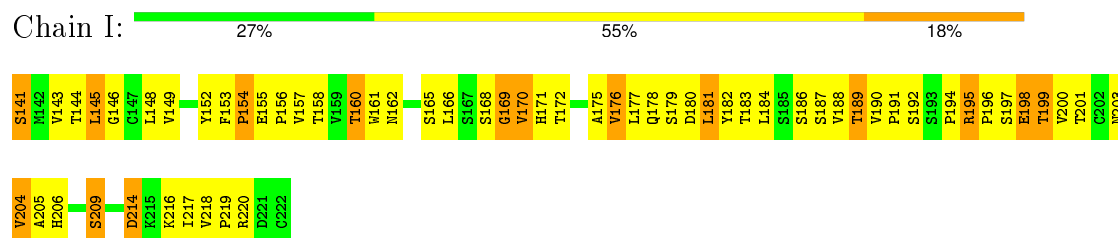
- Molecule 1: PROTEIN (IGG1 ANTIBODY 1696 (light chain))



- Molecule 2: PROTEIN (IGG1 ANTIBODY 1696 (variable heavy chain))



- Molecule 3: PROTEIN (IGG1 ANTIBODY 1696 (constant heavy chain))



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	46.40 Å 58.80 Å 143.00 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 3.00	Depositor
% Data completeness (in resolution range)	96.0 (8.00-3.00)	Depositor
$R_{merge}$	0.13	Depositor
$R_{sym}$	0.13	Depositor
Refinement program	X-PLOR 3.843	Depositor
R, $R_{free}$	0.183 , 0.280	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3355	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	14.0	wwPDB-VP

## 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	L	0.57	0/1727	0.87	1/2344 (0.0%)
2	H	0.66	0/1093	0.90	1/1488 (0.1%)
3	I	0.63	0/621	0.91	1/851 (0.1%)
All	All	0.61	0/3441	0.88	3/4683 (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(°)	Ideal(°)
2	H	98	ASP	N-CA-C	-7.32	91.23	111.00
3	I	177	LEU	CA-CB-CG	6.56	130.38	115.30
1	L	181	LEU	CA-CB-CG	6.23	129.64	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	L	36	TYR	Sidechain

### 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	L	1686	0	1621	118	0
2	H	1063	0	1007	61	0
3	I	606	0	604	58	0
All	All	3355	0	3232	222	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 34.

All (222) 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:4:MET:HB3	1:L:99:GLY:HA2	1.38	1.02
2:H:40:ALA:HB3	2:H:44:ASP:HB3	1.57	0.86
2:H:61:ASP:HA	2:H:64:LYS:HD2	1.59	0.83
2:H:35:HIS:HB2	2:H:93:VAL:HG22	1.62	0.81
1:L:83:LEU:HD13	1:L:104:LEU:HD13	1.64	0.80
2:H:84:ASN:O	2:H:87:THR:HG22	1.81	0.80
3:I:143:VAL:HG23	3:I:192:SER:HA	1.65	0.79
3:I:145:LEU:HD21	3:I:195:ARG:HG3	1.65	0.78
2:H:39:GLN:HA	2:H:44:ASP:O	1.84	0.78
3:I:195:ARG:HB3	3:I:196:PRO:HD3	1.67	0.76
3:I:143:VAL:CG2	3:I:192:SER:HA	2.15	0.75
3:I:168:SER:O	3:I:170:VAL:HG12	1.87	0.74
2:H:18:VAL:HG22	2:H:82(C):LEU:HD11	1.70	0.74
1:L:18:GLN:HG3	1:L:75:ILE:O	1.87	0.73
2:H:35:HIS:O	2:H:92:CYS:HA	1.88	0.73
1:L:27(B):ILE:HG13	1:L:31:THR:HG23	1.71	0.72
1:L:33:LEU:HD22	1:L:71:PHE:CG	2.25	0.72
1:L:193:THR:CG2	1:L:206:VAL:HG13	2.20	0.71
1:L:25:SER:HB3	1:L:69:THR:HG22	1.74	0.69
2:H:33:SER:HB2	2:H:51:ILE:O	1.93	0.69
2:H:61:ASP:O	2:H:64:LYS:HB2	1.94	0.68
1:L:32:TYR:O	1:L:90:GLN:HA	1.93	0.68
3:I:170:VAL:HA	3:I:187:SER:O	1.94	0.68
2:H:54:THR:O	2:H:55:ASP:HB2	1.94	0.68
3:I:162:ASN:ND2	3:I:199:THR:HG22	2.07	0.67
1:L:190:ASN:O	1:L:210:ASN:HA	1.95	0.67
1:L:108:ARG:HD2	1:L:171:SER:HB2	1.76	0.66
1:L:195:GLU:HG2	1:L:206:VAL:HG22	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:93:VAL:HG21	2:H:100(C):PHE:CG	2.31	0.65
2:H:18:VAL:O	2:H:81:GLN:HG3	1.96	0.65
1:L:136:LEU:HD21	1:L:146:VAL:HB	1.79	0.64
1:L:33:LEU:HD22	1:L:71:PHE:CB	2.28	0.63
1:L:146:VAL:HG21	1:L:194:CYS:SG	2.38	0.63
2:H:18:VAL:CG2	2:H:82(C):LEU:HD11	2.27	0.63
1:L:197:THR:OG1	1:L:204:PRO:HB3	1.98	0.63
3:I:158:THR:OG1	3:I:205:ALA:HB3	1.98	0.63
2:H:119:PRO:HB3	3:I:152:TYR:HB3	1.80	0.63
2:H:12:LYS:HE2	2:H:18:VAL:HG13	1.80	0.62
1:L:3:LEU:HD22	1:L:3:LEU:H	1.64	0.62
2:H:69:PHE:HE1	2:H:80:LEU:HD12	1.64	0.62
2:H:55:ASP:O	2:H:57:PRO:HD3	2.00	0.62
3:I:162:ASN:HB3	3:I:165:SER:OG	2.00	0.61
3:I:162:ASN:OD1	3:I:166:LEU:HD11	2.01	0.61
1:L:38:GLN:HE22	2:H:39:GLN:HE22	1.49	0.60
3:I:191:PRO:HB2	3:I:194:PRO:HD2	1.83	0.60
3:I:169:GLY:O	3:I:188:VAL:HA	2.00	0.60
2:H:37:VAL:HG21	2:H:103:TRP:HZ3	1.66	0.60
1:L:181:LEU:HD23	1:L:186:TYR:HB2	1.82	0.60
2:H:12:LYS:CE	2:H:18:VAL:HG13	2.32	0.59
1:L:33:LEU:HD22	1:L:71:PHE:CD1	2.36	0.59
1:L:117:ILE:HD13	1:L:208:SER:HA	1.82	0.59
1:L:183:LYS:O	1:L:187:GLU:HB2	2.02	0.59
2:H:2:ILE:HG21	2:H:94:ARG:NH1	2.18	0.59
1:L:110:ASP:HA	1:L:140:TYR:HD1	1.67	0.59
2:H:51:ILE:HD13	2:H:71:LEU:HD23	1.84	0.59
1:L:181:LEU:HD22	1:L:181:LEU:H	1.67	0.58
1:L:128:GLY:O	1:L:183:LYS:N	2.37	0.58
3:I:195:ARG:NH2	3:I:217:ILE:HG22	2.18	0.58
1:L:33:LEU:HD22	1:L:71:PHE:HB3	1.85	0.58
1:L:83:LEU:HD12	1:L:105:GLU:HA	1.85	0.57
1:L:61:ARG:CZ	1:L:79:GLU:HG3	2.34	0.57
1:L:156:GLN:H	1:L:156:GLN:CD	2.06	0.57
3:I:203:ASN:HA	3:I:214:ASP:HB3	1.85	0.57
1:L:160:LEU:HD21	3:I:178:GLN:HA	1.86	0.57
3:I:162:ASN:HB2	3:I:166:LEU:HG	1.86	0.57
1:L:54:ARG:HG2	1:L:54:ARG:HH11	1.69	0.56
1:L:35:TRP:CE2	1:L:73:LEU:HB2	2.40	0.56
1:L:4:MET:HB3	1:L:99:GLY:CA	2.25	0.56
1:L:27(A):THR:HG22	1:L:27(B):ILE:N	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:155:GLU:HB3	3:I:156:PRO:HA	1.87	0.56
3:I:172:THR:HG23	3:I:184:LEU:HD21	1.87	0.56
2:H:39:GLN:O	2:H:88:ALA:HB1	2.05	0.55
1:L:39:LYS:HE2	1:L:81:GLU:O	2.06	0.55
1:L:106:ILE:H	1:L:166:GLN:HE22	1.54	0.55
1:L:37:LEU:HB2	1:L:47:LEU:HD11	1.89	0.55
2:H:66:ARG:HD2	2:H:82(B):ASN:HB2	1.89	0.55
3:I:191:PRO:HB2	3:I:194:PRO:CD	2.37	0.55
2:H:2:ILE:HG23	2:H:27:TYR:CD1	2.42	0.54
2:H:40:ALA:HB3	2:H:44:ASP:CB	2.33	0.54
2:H:124:LEU:HB2	3:I:146:GLY:C	2.28	0.54
2:H:84:ASN:O	2:H:87:THR:CG2	2.56	0.53
3:I:144:THR:HA	3:I:189:THR:HA	1.90	0.53
1:L:89:PHE:CG	1:L:90:GLN:N	2.76	0.53
2:H:99:TYR:HD1	2:H:99:TYR:N	2.06	0.53
3:I:144:THR:OG1	3:I:189:THR:HG22	2.08	0.53
3:I:157:VAL:HG12	3:I:206:HIS:HB2	1.90	0.53
3:I:194:PRO:O	3:I:197:SER:HB2	2.09	0.53
2:H:126:PRO:HD3	3:I:145:LEU:HD12	1.91	0.52
1:L:146:VAL:CG2	1:L:194:CYS:SG	2.97	0.52
1:L:128:GLY:O	1:L:183:LYS:HB2	2.10	0.52
3:I:171:HIS:O	3:I:186:SER:HA	2.10	0.52
1:L:160:LEU:HD13	3:I:178:GLN:NE2	2.25	0.52
1:L:4:MET:HG3	1:L:23:CYS:SG	2.49	0.51
1:L:12:PRO:HA	1:L:105:GLU:O	2.09	0.51
1:L:166:GLN:HG2	1:L:171:SER:HA	1.91	0.51
1:L:83:LEU:N	1:L:83:LEU:HD22	2.26	0.51
3:I:157:VAL:HA	3:I:205:ALA:O	2.11	0.51
1:L:86:TYR:O	1:L:101:GLY:HA2	2.11	0.51
1:L:38:GLN:HE22	2:H:39:GLN:NE2	2.09	0.50
2:H:2:ILE:HG21	2:H:94:ARG:HH11	1.76	0.50
2:H:37:VAL:HG21	2:H:103:TRP:CZ3	2.44	0.50
1:L:2:VAL:HG22	1:L:27:GLN:HB2	1.94	0.50
1:L:36:TYR:HB3	1:L:44:PRO:HB3	1.94	0.50
3:I:162:ASN:OD1	3:I:166:LEU:CD1	2.59	0.50
2:H:99:TYR:CD1	2:H:99:TYR:N	2.78	0.50
1:L:150:ILE:HG21	1:L:189:HIS:CD2	2.47	0.49
1:L:167:ASP:N	1:L:172:THR:O	2.43	0.49
2:H:60:ALA:O	2:H:62:ASP:N	2.46	0.49
2:H:2:ILE:HG23	2:H:27:TYR:HD1	1.75	0.49
3:I:176:VAL:O	3:I:182:TYR:CD1	2.65	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:148:TRP:HZ2	1:L:177:SER:O	1.95	0.49
3:I:143:VAL:O	3:I:189:THR:HA	2.13	0.49
3:I:190:VAL:HG21	3:I:195:ARG:HB2	1.95	0.49
1:L:120:PRO:HD3	1:L:132:VAL:HG22	1.95	0.49
1:L:150:ILE:N	1:L:153:SER:O	2.46	0.49
1:L:189:HIS:HB2	1:L:192:TYR:OH	2.13	0.48
2:H:119:PRO:CB	3:I:152:TYR:HB3	2.42	0.48
1:L:121:SER:O	1:L:125:LEU:HB2	2.13	0.48
2:H:93:VAL:HG23	2:H:94:ARG:N	2.27	0.48
1:L:181:LEU:CB	1:L:185:GLU:HB2	2.43	0.48
1:L:147:LYS:NZ	1:L:147:LYS:HB3	2.28	0.48
3:I:141:SER:HB3	3:I:192:SER:OG	2.14	0.48
1:L:27(B):ILE:HD11	1:L:71:PHE:CE2	2.48	0.48
1:L:181:LEU:HB3	1:L:185:GLU:HB2	1.95	0.48
1:L:193:THR:HG21	1:L:206:VAL:HG13	1.92	0.48
1:L:160:LEU:HD13	3:I:178:GLN:HE22	1.79	0.48
3:I:190:VAL:CG2	3:I:195:ARG:HB2	2.43	0.48
1:L:146:VAL:HG11	1:L:177:SER:HB2	1.96	0.47
3:I:175:ALA:HA	3:I:183:THR:O	2.14	0.47
2:H:14:PRO:HA	2:H:82(C):LEU:O	2.14	0.47
1:L:146:VAL:HG11	1:L:177:SER:CB	2.44	0.47
1:L:38:GLN:NE2	2:H:39:GLN:HE22	2.12	0.47
1:L:164:THR:HG22	1:L:165:ASP:O	2.14	0.47
1:L:159:VAL:HA	1:L:178:THR:O	2.15	0.47
1:L:85:ILE:HG12	1:L:103:LYS:HD2	1.96	0.47
1:L:52:SER:HB3	1:L:64:GLY:O	2.14	0.47
1:L:37:LEU:HD21	1:L:39:LYS:HE3	1.97	0.47
3:I:153:PHE:CD2	3:I:154:PRO:HA	2.50	0.47
1:L:96:PRO:HD2	2:H:47:TYR:CD2	2.50	0.47
1:L:148:TRP:HE1	1:L:177:SER:HG	1.62	0.46
2:H:56:GLU:HA	2:H:57:PRO:HD2	1.87	0.46
2:H:13:LYS:O	2:H:16:GLU:HG3	2.15	0.46
1:L:113:PRO:HB3	1:L:136:LEU:O	2.15	0.46
3:I:180:ASP:C	3:I:181:LEU:HD22	2.35	0.46
1:L:113:PRO:HG3	1:L:139:PHE:HB3	1.98	0.46
1:L:27(D):HIS:CG	1:L:27(E):ASN:N	2.83	0.46
2:H:2:ILE:HG12	2:H:94:ARG:HH12	1.81	0.46
3:I:158:THR:O	3:I:204:VAL:HA	2.16	0.46
2:H:20:ILE:O	2:H:79:PHE:HA	2.16	0.46
1:L:108:ARG:NH1	1:L:109:ALA:O	2.48	0.45
1:L:11:LEU:HD12	1:L:104:LEU:CD2	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:83:LEU:CD2	1:L:83:LEU:H	2.28	0.45
1:L:209:PHE:HD1	1:L:210:ASN:O	2.00	0.45
1:L:83:LEU:N	1:L:83:LEU:CD2	2.80	0.45
1:L:39:LYS:HB2	1:L:42:GLN:CD	2.36	0.45
1:L:7:THR:HA	1:L:8:PRO:HA	1.76	0.45
3:I:192:SER:O	3:I:196:PRO:HD2	2.17	0.45
1:L:25:SER:HB3	1:L:27:GLN:O	2.17	0.45
2:H:87:THR:O	2:H:88:ALA:HB2	2.17	0.45
1:L:27(A):THR:CG2	1:L:27(C):VAL:HG12	2.46	0.45
2:H:6:GLN:HE21	2:H:6:GLN:HB3	1.53	0.45
1:L:11:LEU:HD12	1:L:104:LEU:HD23	1.98	0.44
3:I:157:VAL:HG23	3:I:157:VAL:O	2.17	0.44
2:H:122:TYR:HB2	3:I:148:LEU:HB3	1.99	0.44
1:L:133:VAL:HG12	1:L:134:CYS:N	2.32	0.44
1:L:126:THR:O	1:L:128:GLY:N	2.50	0.44
1:L:48:ILE:HA	1:L:54:ARG:HA	1.99	0.44
3:I:155:GLU:HB3	3:I:156:PRO:CA	2.46	0.44
2:H:82(B):ASN:O	2:H:83:LYS:NZ	2.48	0.44
1:L:149:LYS:HA	1:L:153:SER:O	2.18	0.44
1:L:27(D):HIS:HB2	1:L:92:SER:HA	1.99	0.44
3:I:201:THR:HA	3:I:216:LYS:HA	2.00	0.44
1:L:27(D):HIS:ND1	1:L:27(E):ASN:N	2.63	0.44
1:L:148:TRP:CD1	1:L:159:VAL:HG21	2.53	0.43
1:L:89:PHE:CE1	1:L:96:PRO:HB3	2.53	0.43
2:H:119:PRO:CA	3:I:152:TYR:HB3	2.48	0.43
2:H:43:GLY:O	2:H:44:ASP:C	2.56	0.43
1:L:63:SER:O	1:L:73:LEU:HD22	2.18	0.43
1:L:175:MET:HG2	1:L:176:SER:N	2.33	0.43
1:L:31:THR:HG21	1:L:71:PHE:CZ	2.53	0.42
3:I:206:HIS:HD2	3:I:209:SER:OG	2.01	0.42
1:L:46:LEU:HD11	1:L:49:TYR:HB3	2.01	0.42
1:L:15:LEU:N	1:L:15:LEU:HD23	2.35	0.42
1:L:62:PHE:CD1	1:L:62:PHE:N	2.87	0.42
3:I:160:THR:HB	3:I:203:ASN:HB2	2.02	0.42
3:I:218:VAL:HA	3:I:219:PRO:HD3	1.88	0.42
1:L:3:LEU:HD22	1:L:3:LEU:N	2.34	0.42
1:L:192:TYR:O	1:L:208:SER:HA	2.20	0.42
1:L:33:LEU:CD2	1:L:71:PHE:HB3	2.47	0.42
2:H:59:PHE:CE2	2:H:69:PHE:CE2	3.08	0.42
3:I:170:VAL:C	3:I:171:HIS:HD2	2.23	0.42
1:L:140:TYR:C	1:L:140:TYR:CD1	2.92	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:160:THR:HG22	3:I:203:ASN:OD1	2.20	0.42
1:L:199:LYS:HE2	1:L:199:LYS:HB2	1.90	0.41
2:H:6:GLN:NE2	2:H:107:THR:HG22	2.34	0.41
1:L:31:THR:HG21	1:L:71:PHE:CE1	2.54	0.41
3:I:155:GLU:CB	3:I:156:PRO:HA	2.49	0.41
1:L:34:GLU:O	1:L:88:CYS:HA	2.20	0.41
1:L:140:TYR:O	1:L:198:HIS:HE1	2.04	0.41
1:L:160:LEU:CD2	3:I:178:GLN:HA	2.49	0.41
2:H:66:ARG:NH2	2:H:86:ASP:OD2	2.54	0.41
2:H:41:PRO:HD3	2:H:88:ALA:HA	2.03	0.41
1:L:27(B):ILE:O	1:L:31:THR:HA	2.21	0.41
1:L:49:TYR:CZ	2:H:96:ARG:NH1	2.88	0.41
2:H:67:PHE:CE2	2:H:82:ILE:HD13	2.56	0.41
1:L:81:GLU:CD	1:L:81:GLU:N	2.73	0.41
1:L:182:THR:HG1	1:L:184:ASP:CG	2.23	0.41
2:H:48:VAL:HA	2:H:63:PHE:CD1	2.56	0.41
1:L:33:LEU:HD23	1:L:51:VAL:HA	2.02	0.41
2:H:47:TYR:OH	2:H:50:TRP:HE3	2.04	0.41
1:L:49:TYR:O	1:L:53:ASN:HB2	2.20	0.41
1:L:135:PHE:HB3	1:L:137:ASN:OD1	2.21	0.40
3:I:206:HIS:CD2	3:I:209:SER:OG	2.74	0.40
2:H:1:GLN:O	2:H:3:GLN:HG3	2.21	0.40
3:I:149:VAL:HG22	3:I:204:VAL:HG11	2.04	0.40
3:I:160:THR:CG2	3:I:161:TRP:N	2.85	0.40
2:H:83:LYS:HZ2	2:H:83:LYS:HA	1.86	0.40
1:L:141:PRO:HD2	1:L:199:LYS:HD2	2.02	0.40
3:I:145:LEU:HD23	3:I:200:VAL:HG11	2.03	0.40
1:L:95:PRO:HA	1:L:96:PRO:HD3	1.86	0.40
1:L:54:ARG:HG2	1:L:54:ARG:NH1	2.36	0.40
1:L:49:TYR:CD1	1:L:53:ASN:HB2	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	L	214/216 (99%)	184 (86%)	23 (11%)	7 (3%)	5	26
2	H	133/135 (98%)	106 (80%)	17 (13%)	10 (8%)	1	6
3	I	80/82 (98%)	67 (84%)	7 (9%)	6 (8%)	1	6
All	All	427/433 (99%)	357 (84%)	47 (11%)	23 (5%)	2	14

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	137	ASN
2	H	61	ASP
2	H	100(C)	PHE
3	I	179	SER
2	H	44	ASP
2	H	65	GLY
3	I	198	GLU
1	L	67	SER
1	L	188	ARG
2	H	82(C)	LEU
2	H	88	ALA
2	H	116	THR
1	L	78	VAL
2	H	75	THR
3	I	154	PRO
1	L	80	ALA
1	L	127	SER
2	H	82(B)	ASN
3	I	199	THR
1	L	51	VAL
3	I	169	GLY
2	H	100(B)	ILE
3	I	176	VAL

### 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	193/193 (100%)	167 (86%)	26 (14%)	5	20
2	H	116/116 (100%)	92 (79%)	24 (21%)	1	7
3	I	75/75 (100%)	63 (84%)	12 (16%)	3	15
All	All	384/384 (100%)	322 (84%)	62 (16%)	3	14

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	3	LEU
1	L	4	MET
1	L	9	LEU
1	L	13	VAL
1	L	15	LEU
1	L	25	SER
1	L	26	SER
1	L	27(C)	VAL
1	L	30	ASN
1	L	48	ILE
1	L	52	SER
1	L	54	ARG
1	L	74	LYS
1	L	79	GLU
1	L	83	LEU
1	L	90	GLN
1	L	97	THR
1	L	104	LEU
1	L	105	GLU
1	L	143	ASP
1	L	147	LYS
1	L	156	GLN
1	L	171	SER
1	L	181	LEU
1	L	187	GLU
1	L	199	LYS
2	H	1	GLN
2	H	6	GLN
2	H	14	PRO
2	H	18	VAL
2	H	21	SER
2	H	26	ASN

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Mol	Chain	Res	Type
2	H	31	ASP
2	H	41	PRO
2	H	55	ASP
2	H	80	LEU
2	H	82(B)	ASN
2	H	83	LYS
2	H	84	ASN
2	H	87	THR
2	H	93	VAL
2	H	94	ARG
2	H	96	ARG
2	H	98	ASP
2	H	105	GLN
2	H	107	THR
2	H	108	THR
2	H	120	SER
2	H	124	LEU
2	H	128	SER
3	I	141	SER
3	I	145	LEU
3	I	160	THR
3	I	170	VAL
3	I	181	LEU
3	I	189	THR
3	I	195	ARG
3	I	198	GLU
3	I	204	VAL
3	I	209	SER
3	I	214	ASP
3	I	220	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	28	ASN
1	L	38	GLN
1	L	42	GLN
1	L	93	HIS
1	L	157	ASN
1	L	166	GLN
1	L	190	ASN
2	H	39	GLN

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Mol	Chain	Res	Type
3	I	171	HIS
3	I	178	GLN
3	I	206	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 [i](#)

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

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.