Full wwPDB X-ray Structure Validation Report (i)

May 29, 2024 – 06:49 PM EDT

PDB ID	:	1GHF
Title	:	ANTI-ANTI-IDIOTYPE GH1002 FAB FRAGMENT
Authors	:	Ban, N.; Day, J.; Wang, X.; Ferrone, S.; McPherson, A.
Deposited on	:	1995-11-30
Resolution	:	2.70 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*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

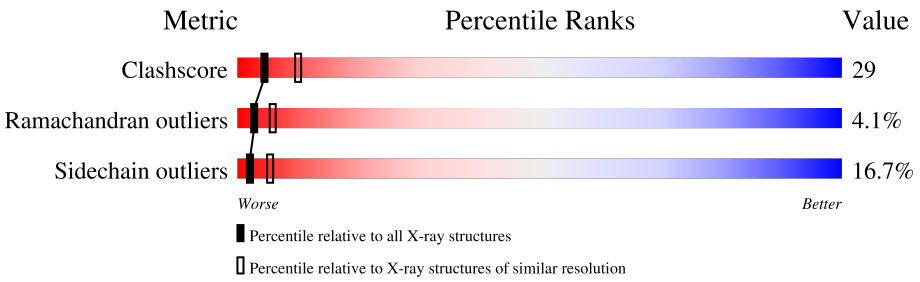
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)

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 >=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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	211	46% 45% 7% •
2	H	212	48% 39% 13%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3273 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 ANTI-ANTI-IDIOTYPE GH1002 FAB FRAGMENT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	L	211	1646	1018	278	344	6	0	0	0

- Molecule 2 is a protein called ANTI-ANTI-IDIOTYPE GH1002 FAB FRAGMENT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	H	212	1627	1041	263	315	8	44	0	0

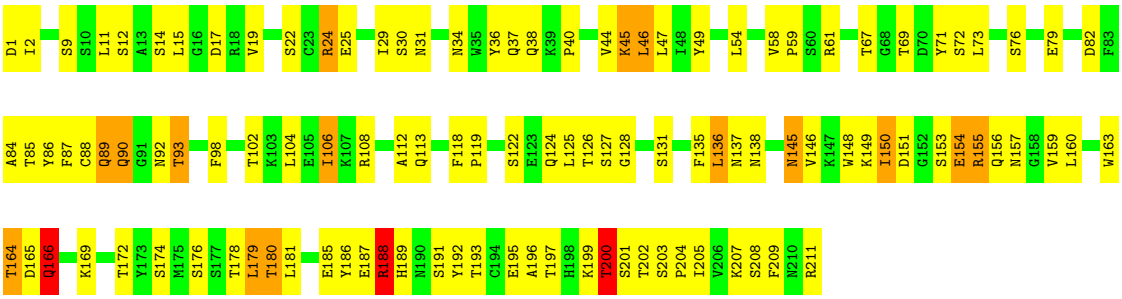
3 Residue-property plots (i)

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. 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. 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.

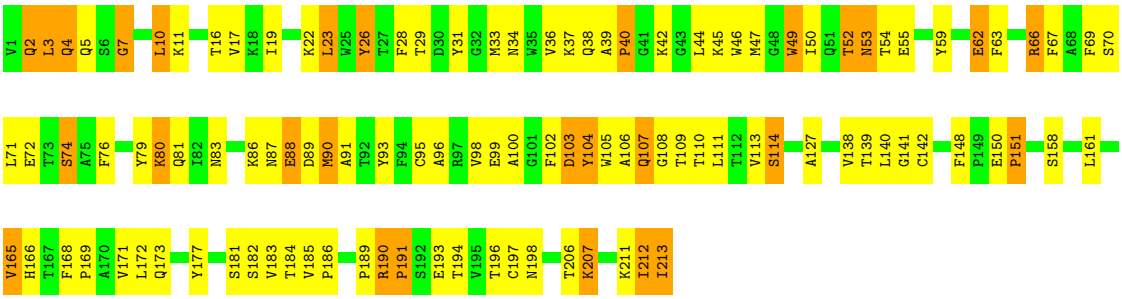
• Molecule 1: ANTI-ANTI-IDIOTYPE GH1002 FAB FRAGMENT

Chain L: 46% 45% 7%



• Molecule 2: ANTI-ANTI-IDIOTYPE GH1002 FAB FRAGMENT

Chain H: 48% 39% 13%



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α, β, γ	150.40Å 150.40Å 43.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.70	Depositor
% Data completeness (in resolution range)	(Not available) (30.00-2.70)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.210 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3273	wwPDB-VP
Average B, all atoms (Å^2)	10.0	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	0.80	0/1679	1.01	3/2276 (0.1%)
2	H	0.78	1/1673 (0.1%)	1.01	6/2283 (0.3%)
All	All	0.79	1/3352 (0.0%)	1.01	9/4559 (0.2%)

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
2	H	0	3
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	62	GLU	CG-CD	5.17	1.59	1.51

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	L	73	LEU	CA-CB-CG	7.15	131.74	115.30
2	H	111	LEU	CA-CB-CG	7.07	131.56	115.30
2	H	66	ARG	NE-CZ-NH2	-6.77	116.92	120.30
2	H	49	TRP	CB-CA-C	-6.43	97.55	110.40
1	L	76	SER	N-CA-C	6.05	127.33	111.00
2	H	100	ALA	N-CA-C	-6.03	94.72	111.00
2	H	23	LEU	N-CA-C	5.43	125.67	111.00
2	H	114	SER	N-CA-C	5.43	125.67	111.00
1	L	209	PHE	N-CA-C	-5.10	97.23	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	104	TYR	Sidechain
2	H	26	TYR	Sidechain
2	H	49	TRP	Mainchain
1	L	71	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	1646	0	1569	88	0
2	H	1627	0	1594	98	0
All	All	3273	0	3163	182	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 29.

All (182) 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:H:16:THR:HG22	2:H:83:ASN:HA	1.37	1.04
2:H:70:SER:HB2	2:H:79:TYR:HB2	1.39	1.02
2:H:5:GLN:NE2	2:H:95:CYS:SG	2.35	0.99
2:H:80:LYS:NZ	2:H:80:LYS:HB3	1.85	0.90
2:H:7:GLY:HA3	2:H:109:THR:HG23	1.55	0.88
2:H:80:LYS:HB3	2:H:80:LYS:HZ2	1.37	0.84
1:L:148:TRP:CZ3	1:L:179:LEU:HB2	2.13	0.84
2:H:33:MET:HE1	2:H:95:CYS:HB2	1.64	0.79
2:H:186:PRO:O	2:H:189:PRO:HD2	1.83	0.79
1:L:14:SER:O	1:L:17:ASP:HB2	1.83	0.78
2:H:138:VAL:HG12	2:H:185:VAL:O	1.83	0.77
1:L:148:TRP:O	1:L:154:GLU:HA	1.85	0.77
1:L:9:SER:O	1:L:102:THR:HA	1.88	0.74
2:H:19:ILE:HB	2:H:80:LYS:NZ	2.03	0.73
1:L:25:GLU:OE1	1:L:29:ILE:HB	1.88	0.73
2:H:2:GLN:NE2	2:H:3:LEU:HD12	2.03	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:188:ARG:HH11	1:L:188:ARG:HG3	1.54	0.72
2:H:47:MET:HA	2:H:63:PHE:CD2	2.24	0.72
2:H:190:ARG:HB3	2:H:191:PRO:HD3	1.72	0.72
2:H:42:LYS:HB3	2:H:42:LYS:NZ	2.05	0.71
1:L:148:TRP:HE1	1:L:159:VAL:HG11	1.56	0.70
2:H:5:GLN:HG2	2:H:106:ALA:HB1	1.72	0.70
2:H:3:LEU:HD23	2:H:23:LEU:HD21	1.73	0.69
1:L:159:VAL:O	1:L:160:LEU:HD13	1.92	0.69
2:H:107:GLN:HE21	2:H:107:GLN:H	1.41	0.68
2:H:23:LEU:HD13	2:H:28:PHE:HE1	1.60	0.67
1:L:38:GLN:O	1:L:84:ALA:HB1	1.96	0.66
1:L:113:GLN:HG3	1:L:136:LEU:HB3	1.77	0.66
2:H:11:LYS:O	2:H:113:VAL:HA	1.94	0.66
1:L:149:LYS:HA	1:L:153:SER:O	1.95	0.66
1:L:36:TYR:OH	1:L:89:GLN:NE2	2.29	0.65
1:L:131:SER:HA	1:L:180:THR:HA	1.78	0.65
1:L:90:GLN:HG2	1:L:92:ASN:H	1.61	0.65
2:H:3:LEU:CD2	2:H:23:LEU:HD21	2.27	0.65
2:H:28:PHE:HE2	2:H:71:LEU:HD11	1.62	0.65
2:H:34:ASN:HD22	2:H:46:TRP:HE1	1.43	0.64
2:H:55:GLU:HG2	2:H:71:LEU:HD23	1.79	0.64
2:H:33:MET:CE	2:H:95:CYS:HB2	2.28	0.63
1:L:155:ARG:HD2	1:L:155:ARG:N	2.14	0.63
2:H:39:ALA:HB1	2:H:40:PRO:HD2	1.81	0.63
2:H:19:ILE:HB	2:H:80:LYS:HZ3	1.64	0.62
1:L:149:LYS:O	1:L:151:ASP:N	2.33	0.62
2:H:96:ALA:HA	2:H:104:TYR:O	2.00	0.62
2:H:213:ILE:H	2:H:213:ILE:HD13	1.63	0.61
1:L:29:ILE:HG22	1:L:29:ILE:O	1.99	0.61
1:L:131:SER:OG	1:L:180:THR:HB	2.01	0.60
1:L:148:TRP:CD1	1:L:159:VAL:HG21	2.38	0.59
1:L:136:LEU:C	1:L:137:ASN:HD22	2.06	0.59
2:H:38:GLN:HB2	2:H:44:LEU:HD23	1.84	0.59
1:L:29:ILE:HD12	1:L:90:GLN:HG3	1.85	0.59
2:H:3:LEU:HD12	2:H:3:LEU:H	1.68	0.59
2:H:7:GLY:HA3	2:H:109:THR:CG2	2.29	0.58
1:L:106:ILE:H	1:L:166:GLN:HE22	1.50	0.58
2:H:189:PRO:O	2:H:193:GLU:HB2	2.03	0.58
2:H:37:LYS:O	2:H:45:LYS:HG2	2.03	0.58
2:H:39:ALA:O	2:H:42:LYS:HB2	2.02	0.58
1:L:153:SER:HB3	1:L:155:ARG:HE	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:23:LEU:CD2	2:H:26:TYR:HE1	2.17	0.57
1:L:36:TYR:CD1	1:L:46:LEU:HA	2.39	0.56
2:H:2:GLN:NE2	2:H:3:LEU:H	2.02	0.56
2:H:5:GLN:HG3	2:H:108:GLY:H	1.70	0.56
2:H:3:LEU:HD12	2:H:3:LEU:N	2.21	0.56
2:H:66:ARG:NH1	2:H:86:LYS:HE3	2.21	0.56
2:H:38:GLN:O	2:H:91:ALA:HB1	2.06	0.56
1:L:153:SER:HB3	1:L:155:ARG:NE	2.21	0.56
1:L:150:ILE:HG12	1:L:153:SER:HB2	1.87	0.55
1:L:160:LEU:HD21	2:H:173:GLN:OE1	2.06	0.55
1:L:138:ASN:HA	1:L:172:THR:OG1	2.05	0.55
2:H:34:ASN:ND2	2:H:46:TRP:HE1	2.02	0.55
2:H:3:LEU:HD13	2:H:105:TRP:O	2.07	0.55
2:H:16:THR:CG2	2:H:83:ASN:HA	2.24	0.55
1:L:24:ARG:HA	1:L:69:THR:O	2.07	0.54
1:L:44:VAL:HG12	1:L:45:LYS:N	2.22	0.54
1:L:148:TRP:NE1	1:L:159:VAL:HG21	2.23	0.54
1:L:155:ARG:O	1:L:159:VAL:HG23	2.07	0.54
1:L:160:LEU:HD12	2:H:171:VAL:HG21	1.90	0.53
1:L:188:ARG:HG3	1:L:188:ARG:NH1	2.22	0.53
1:L:150:ILE:HG22	1:L:192:TYR:HA	1.91	0.52
1:L:196:ALA:O	1:L:205:ILE:HG22	2.08	0.52
1:L:211:ARG:HH11	1:L:211:ARG:HG3	1.74	0.52
2:H:5:GLN:H	2:H:107:GLN:NE2	2.09	0.51
2:H:165:VAL:HA	2:H:182:SER:O	2.11	0.51
1:L:34:ASN:OD1	1:L:49:TYR:HA	2.10	0.51
1:L:150:ILE:HA	1:L:193:THR:HG23	1.93	0.51
1:L:164:THR:HG23	1:L:165:ASP:N	2.25	0.51
1:L:29:ILE:O	1:L:29:ILE:CG2	2.59	0.50
2:H:23:LEU:HD23	2:H:26:TYR:HE1	1.76	0.50
2:H:190:ARG:HG2	2:H:190:ARG:HH11	1.76	0.50
2:H:29:THR:HA	2:H:52:THR:HB	1.93	0.50
1:L:178:THR:HG22	1:L:180:THR:HG22	1.94	0.49
1:L:150:ILE:CG1	1:L:153:SER:HB2	2.43	0.49
1:L:25:GLU:OE2	1:L:29:ILE:HG21	2.12	0.49
2:H:22:LYS:C	2:H:23:LEU:HD12	2.31	0.49
1:L:25:GLU:CD	1:L:29:ILE:HB	2.31	0.49
1:L:36:TYR:HD1	1:L:46:LEU:HA	1.77	0.49
1:L:146:VAL:HB	1:L:195:GLU:O	2.13	0.49
2:H:39:ALA:CB	2:H:40:PRO:HD2	2.41	0.49
2:H:42:LYS:HB3	2:H:42:LYS:HZ2	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:34:ASN:O	2:H:95:CYS:HA	2.13	0.49
1:L:118:PHE:HE2	1:L:135:PHE:HD2	1.61	0.48
1:L:192:TYR:CD1	1:L:192:TYR:N	2.82	0.48
2:H:33:MET:HE3	2:H:34:ASN:N	2.28	0.48
2:H:19:ILE:HD12	2:H:80:LYS:NZ	2.29	0.48
2:H:166:HIS:O	2:H:181:SER:HA	2.14	0.48
1:L:124:GLN:O	1:L:127:SER:HB2	2.13	0.48
1:L:187:GLU:HA	1:L:187:GLU:OE1	2.14	0.48
2:H:168:PHE:HA	2:H:169:PRO:HD3	1.72	0.48
2:H:140:LEU:HD12	2:H:212:ILE:HG21	1.96	0.48
2:H:36:VAL:HG11	2:H:44:LEU:HD22	1.95	0.48
2:H:190:ARG:HG2	2:H:190:ARG:NH1	2.29	0.47
1:L:145:ASN:OD1	1:L:197:THR:HB	2.14	0.47
2:H:47:MET:HA	2:H:63:PHE:CE2	2.50	0.47
1:L:30:SER:OG	1:L:31:ASN:N	2.46	0.46
2:H:66:ARG:NH2	2:H:89:ASP:OD2	2.48	0.46
1:L:11:LEU:HD23	1:L:12:SER:N	2.30	0.46
1:L:59:PRO:HB2	1:L:61:ARG:HG3	1.98	0.46
2:H:90:MET:SD	2:H:113:VAL:HB	2.55	0.46
1:L:44:VAL:HG12	1:L:45:LYS:H	1.81	0.46
2:H:4:GLN:HA	2:H:107:GLN:HE22	1.81	0.46
2:H:107:GLN:HE21	2:H:107:GLN:N	2.10	0.46
1:L:207:LYS:HD2	1:L:207:LYS:HA	1.80	0.45
2:H:17:VAL:O	2:H:81:GLN:HB2	2.16	0.45
2:H:23:LEU:HD13	2:H:28:PHE:CE1	2.47	0.45
2:H:161:LEU:HA	2:H:161:LEU:HD23	1.49	0.45
2:H:28:PHE:CE2	2:H:71:LEU:HD11	2.45	0.45
2:H:140:LEU:O	2:H:182:SER:HA	2.16	0.45
2:H:59:TYR:CE1	2:H:69:PHE:CE2	3.05	0.45
1:L:37:GLN:HB2	1:L:47:LEU:HD11	1.99	0.45
1:L:155:ARG:HG2	1:L:155:ARG:HH11	1.82	0.45
1:L:199:LYS:O	1:L:201:SER:N	2.49	0.45
2:H:93:TYR:O	2:H:108:GLY:HA2	2.17	0.45
1:L:185:GLU:O	1:L:189:HIS:ND1	2.50	0.44
1:L:2:ILE:HD13	1:L:29:ILE:HD11	1.99	0.44
2:H:196:THR:HG22	2:H:211:LYS:HA	1.98	0.44
1:L:122:SER:O	1:L:125:LEU:HB2	2.18	0.44
1:L:138:ASN:OD1	2:H:166:HIS:CE1	2.70	0.44
2:H:23:LEU:HD12	2:H:23:LEU:N	2.32	0.44
1:L:163:TRP:O	2:H:169:PRO:HG2	2.17	0.44
2:H:190:ARG:HD3	2:H:191:PRO:N	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:29:ILE:O	1:L:30:SER:C	2.56	0.44
1:L:211:ARG:HG3	1:L:211:ARG:NH1	2.33	0.44
1:L:113:GLN:H	1:L:113:GLN:CD	2.21	0.43
2:H:107:GLN:HG2	2:H:108:GLY:O	2.18	0.43
2:H:172:LEU:HD13	2:H:177:TYR:CD1	2.53	0.43
1:L:189:HIS:HB2	1:L:192:TYR:OH	2.19	0.43
2:H:3:LEU:HD11	2:H:104:TYR:CB	2.49	0.43
2:H:139:THR:HG23	2:H:183:VAL:O	2.19	0.43
2:H:5:GLN:HG3	2:H:108:GLY:N	2.34	0.43
1:L:112:ALA:HB2	1:L:200:THR:OG1	2.18	0.42
2:H:138:VAL:HG11	2:H:190:ARG:HB2	2.00	0.42
1:L:125:LEU:HD23	1:L:125:LEU:HA	1.88	0.42
2:H:10:LEU:HD11	2:H:148:PHE:CZ	2.54	0.42
1:L:24:ARG:O	1:L:24:ARG:HG2	2.18	0.42
1:L:38:GLN:HB2	1:L:87:PHE:HE2	1.84	0.42
1:L:92:ASN:O	1:L:93:THR:HG23	2.20	0.42
1:L:148:TRP:CH2	1:L:179:LEU:HB2	2.54	0.42
2:H:150:GLU:HA	2:H:151:PRO:HA	1.67	0.42
1:L:106:ILE:N	1:L:166:GLN:HE22	2.17	0.42
2:H:53:ASN:HD22	2:H:54:THR:HG23	1.85	0.42
1:L:193:THR:HA	1:L:208:SER:OG	2.19	0.41
2:H:67:PHE:CD1	2:H:67:PHE:N	2.88	0.41
1:L:90:GLN:HE21	1:L:90:GLN:HB3	1.69	0.41
2:H:190:ARG:CB	2:H:191:PRO:HD3	2.45	0.41
1:L:137:ASN:ND2	1:L:174:SER:HB3	2.35	0.41
1:L:186:TYR:HA	1:L:192:TYR:OH	2.21	0.41
2:H:87:ASN:O	2:H:89:ASP:N	2.53	0.41
2:H:102:PHE:N	2:H:102:PHE:CD1	2.88	0.41
1:L:118:PHE:HA	1:L:119:PRO:HD3	1.70	0.41
2:H:172:LEU:HD12	2:H:172:LEU:HA	1.78	0.41
1:L:88:CYS:O	1:L:98:PHE:HA	2.20	0.41
2:H:31:TYR:C	2:H:52:THR:HG1	2.24	0.41
1:L:86:TYR:CD1	1:L:86:TYR:N	2.89	0.41
2:H:33:MET:HB2	2:H:50:ILE:CG2	2.51	0.41
2:H:53:ASN:HD22	2:H:54:THR:N	2.19	0.41
2:H:127:ALA:HB3	2:H:213:ILE:HG13	2.03	0.41
1:L:15:LEU:C	1:L:17:ASP:H	2.24	0.40
1:L:202:THR:O	1:L:204:PRO:HD3	2.20	0.40
2:H:141:GLY:HA2	2:H:181:SER:O	2.21	0.40
1:L:148:TRP:NE1	1:L:159:VAL:HG11	2.31	0.40
1:L:58:VAL:HA	1:L:59:PRO:HD3	1.95	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:150:ILE:HD11	1:L:155:ARG:NH1	2.37	0.40
2:H:206:THR:C	2:H:207:LYS:HG3	2.42	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	209/211 (99%)	172 (82%)	28 (13%)	9 (4%)	2 5
2	H	210/212 (99%)	171 (81%)	31 (15%)	8 (4%)	3 7
All	All	419/423 (99%)	343 (82%)	59 (14%)	17 (4%)	3 6

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	150	ILE
1	L	200	THR
2	H	40	PRO
2	H	88	GLU
1	L	166	GLN
1	L	188	ARG
2	H	7	GLY
2	H	103	ASP
2	H	191	PRO
1	L	191	SER
2	H	76	PHE
1	L	40	PRO
2	H	74	SER
1	L	82	ASP
1	L	126	THR
1	L	128	GLY

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Mol	Chain	Res	Type
2	H	151	PRO

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	L	190/190 (100%)	157 (83%)	33 (17%)	2 5
2	H	181/181 (100%)	152 (84%)	29 (16%)	2 6
All	All	371/371 (100%)	309 (83%)	62 (17%)	2 5

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

Mol	Chain	Res	Type
1	L	1	ASP
1	L	19	VAL
1	L	22	SER
1	L	24	ARG
1	L	45	LYS
1	L	46	LEU
1	L	54	LEU
1	L	67	THR
1	L	72	SER
1	L	79	GLU
1	L	85	THR
1	L	89	GLN
1	L	90	GLN
1	L	93	THR
1	L	104	LEU
1	L	106	ILE
1	L	108	ARG
1	L	136	LEU
1	L	145	ASN
1	L	154	GLU
1	L	155	ARG
1	L	156	GLN

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Mol	Chain	Res	Type
1	L	157	ASN
1	L	164	THR
1	L	166	GLN
1	L	169	LYS
1	L	176	SER
1	L	179	LEU
1	L	180	THR
1	L	181	LEU
1	L	188	ARG
1	L	200	THR
1	L	203	SER
2	H	2	GLN
2	H	3	LEU
2	H	4	GLN
2	H	10	LEU
2	H	52	THR
2	H	53	ASN
2	H	62	GLU
2	H	72	GLU
2	H	74	SER
2	H	80	LYS
2	H	88	GLU
2	H	90	MET
2	H	98	VAL
2	H	99	GLU
2	H	103	ASP
2	H	107	GLN
2	H	110	THR
2	H	114	SER
2	H	142	CYS
2	H	158	SER
2	H	165	VAL
2	H	184	THR
2	H	190	ARG
2	H	194	THR
2	H	197	CYS
2	H	198	ASN
2	H	207	LYS
2	H	212	ILE
2	H	213	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	6	GLN
1	L	37	GLN
1	L	89	GLN
1	L	90	GLN
1	L	124	GLN
1	L	137	ASN
1	L	161	ASN
1	L	166	GLN
1	L	190	ASN
1	L	210	ASN
2	H	2	GLN
2	H	53	ASN
2	H	107	GLN
2	H	198	ASN

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 monosaccharides 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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.