



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 04:46 pm BST

PDB ID : 5IHZ
Title : Crystal structure of anti-gliadin 1002-1E01 Fab fragment
Authors : Snir, O.; Chen, X.; Gidoni, M.; du Pre, M.F.; Zhao, Y.; Steinsbo, O.; Lundin, K.E.; Yaari, G.; Sollid, L.M.
Deposited on : 2016-03-01
Resolution : 1.64 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 ⓘ symbol.

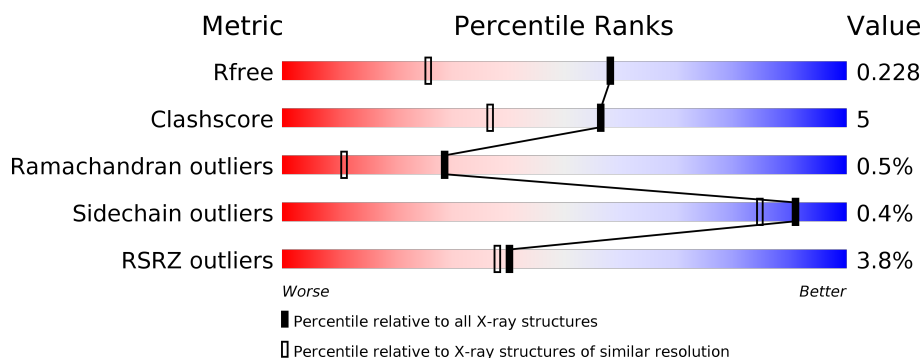
The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

i

X-RAY DIFFRACTION

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}	130704	3122 (1.66-1.62)
Clashscore	141614	3268 (1.66-1.62)
Ramachandran outliers	138981	3215 (1.66-1.62)
Sidechain outliers	138945	3215 (1.66-1.62)
RSRZ outliers	127900	3079 (1.66-1.62)

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 ≥ 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	227	<div> <div></div> <div>86%</div> <div>5%</div> <div>8%</div> </div>
1	C	227	<div> <div></div> <div>83%</div> <div>8%</div> <div>8%</div> </div>
1	E	227	<div> <div></div> <div>86%</div> <div>5%</div> <div>8%</div> </div>
1	H	227	<div> <div></div> <div>86%</div> <div>7%</div> <div>7%</div> </div>
2	B	216	<div> <div></div> <div>88%</div> <div>10%</div> <div>••</div> </div>
2	D	216	<div> <div></div> <div>88%</div> <div>10%</div> <div>•</div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	F	216	<div><div></div><div>6%</div><div>81%</div><div>16%</div><div>••</div></div>
2	L	216	<div><div></div><div>6%</div><div>88%</div><div>11%</div><div>•</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13783 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 1E01 Fab fragment heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	0	0	0
			1543	974	259	304	6			
1	C	208	Total	C	N	O	S	0	0	0
			1546	976	259	305	6			
1	E	209	Total	C	N	O	S	0	0	0
			1547	976	260	305	6			
1	H	211	Total	C	N	O	S	0	0	0
			1565	987	263	308	7			

- Molecule 2 is a protein called 1E01 Fab fragment light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	213	Total	C	N	O	S	0	0	0
			1593	993	270	325	5			
2	D	213	Total	C	N	O	S	0	0	0
			1593	993	270	325	5			
2	F	213	Total	C	N	O	S	0	0	0
			1593	993	270	325	5			
2	L	215	Total	C	N	O	S	0	0	0
			1609	1002	272	330	5			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	165	Total	O	0	0
			165	165		
3	B	105	Total	O	0	0
			105	105		
3	C	146	Total	O	0	0
			146	146		
3	D	133	Total	O	0	0
			133	133		

Continued on next page...

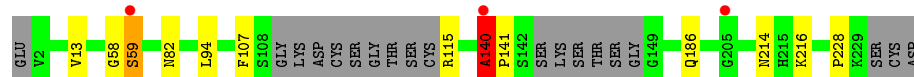
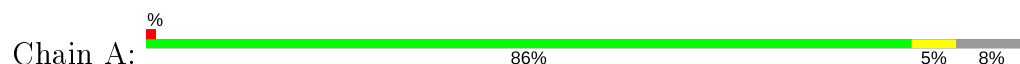
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	196	Total 196	O 196	0	0
3	F	106	Total 106	O 106	0	0
3	H	187	Total 187	O 187	0	0
3	L	156	Total 156	O 156	0	0

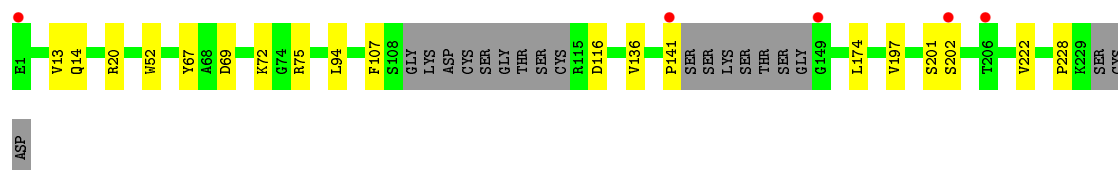
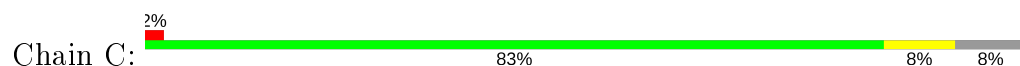
3 Residue-property plots [i](#)

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

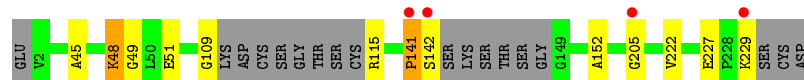
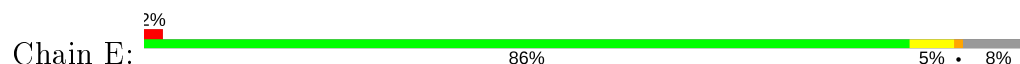
- Molecule 1: 1E01 Fab fragment heavy chain



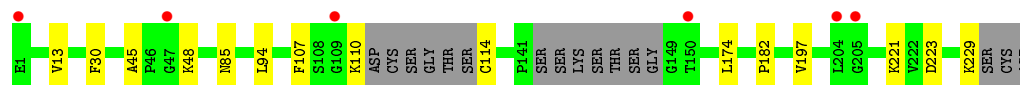
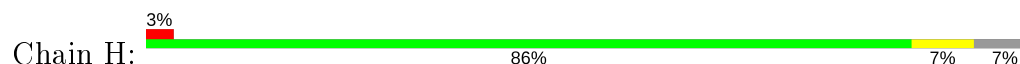
- Molecule 1: 1E01 Fab fragment heavy chain



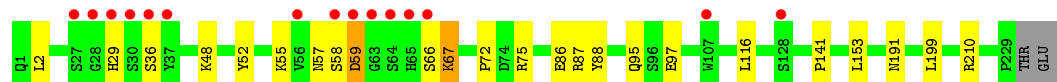
- Molecule 1: 1E01 Fab fragment heavy chain




- Molecule 1: 1E01 Fab fragment heavy chain

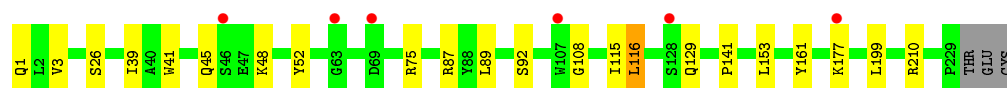


- Molecule 2: 1E01 Fab fragment light chain




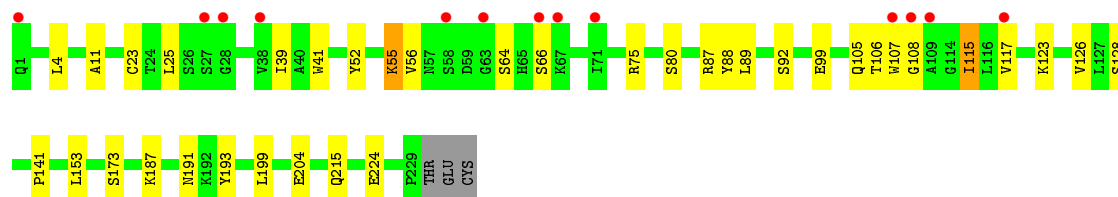
- Molecule 2: 1E01 Fab fragment light chain

Chain D:  3% 88% 10% .




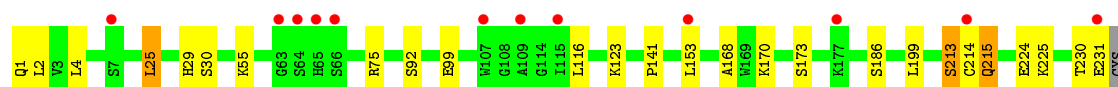
- Molecule 2: 1E01 Fab fragment light chain

Chain F:  6% 81% 16% ..



- Molecule 2: 1E01 Fab fragment light chain

Chain L:  6% 88% 11% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.27Å 147.81Å 95.23Å 90.00° 103.50° 90.00°	Depositor
Resolution (Å)	78.47 – 1.64 78.47 – 1.64	Depositor EDS
% Data completeness (in resolution range)	98.7 (78.47-1.64) 98.7 (78.47-1.64)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 1.64Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155)	Depositor
R, R_{free}	0.200 , 0.229 0.201 , 0.228	Depositor DCC
R_{free} test set	11892 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.317	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13783	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.55% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 ⓘ

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	0.55	0/1578	0.74	2/2145 (0.1%)
1	C	0.53	0/1581	0.65	0/2149
1	E	0.56	0/1582	0.71	1/2150 (0.0%)
1	H	0.55	0/1600	0.73	0/2173
2	B	0.45	0/1630	0.64	0/2219
2	D	0.49	0/1630	0.68	1/2219 (0.0%)
2	F	0.53	1/1630 (0.1%)	0.80	3/2219 (0.1%)
2	L	0.54	0/1646	0.73	1/2241 (0.0%)
All	All	0.52	1/12877 (0.0%)	0.71	8/17515 (0.0%)

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	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	173	SER	C-N	5.11	1.45	1.34

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	115	ILE	CG1-CB-CG2	-11.10	86.98	111.40
1	A	140	ALA	C-N-CD	-10.86	96.72	120.60
2	L	25	LEU	CA-CB-CG	8.31	134.42	115.30
1	A	140	ALA	C-N-CA	7.33	152.79	122.00
2	F	55	LYS	CD-CE-NZ	-6.36	97.08	111.70
2	D	116	LEU	CB-CG-CD2	-5.99	100.81	111.00
2	F	115	ILE	CA-CB-CG1	5.74	121.90	111.00

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	48	LYS	CG-CD-CE	-5.62	95.03	111.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	59	SER	Peptide

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	A	1543	0	1510	14	0
1	C	1546	0	1514	14	0
1	E	1547	0	1513	16	0
1	H	1565	0	1535	10	0
2	B	1593	0	1546	18	0
2	D	1593	0	1546	15	0
2	F	1593	0	1546	30	0
2	L	1609	0	1559	18	0
3	A	165	0	0	4	0
3	B	105	0	0	2	0
3	C	146	0	0	0	0
3	D	133	0	0	0	0
3	E	196	0	0	2	0
3	F	106	0	0	12	0
3	H	187	0	0	0	0
3	L	156	0	0	0	0
All	All	13783	0	12269	125	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 5.

All (125) 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:A:59:SER:HA	1:A:82:ASN:HD21	1.20	1.03
2:F:23:CYS:SG	3:F:403:HOH:O	2.20	0.98
1:A:59:SER:HA	1:A:82:ASN:ND2	1.90	0.87
2:L:170:LYS:HE3	2:L:173:SER:HA	1.57	0.87
2:F:55:LYS:NZ	3:F:302:HOH:O	2.07	0.86
2:D:45:GLN:OE1	2:D:48:LYS:HE3	1.80	0.82
1:E:51:GLU:HB2	3:E:301:HOH:O	1.81	0.80
1:E:45:ALA:HB3	1:E:48:LYS:HE2	1.65	0.79
1:E:227:GLU:O	1:E:229:LYS:NZ	2.17	0.78
2:B:95:GLN:NE2	2:B:97:GLU:OE1	2.16	0.77
1:E:48:LYS:NZ	3:E:301:HOH:O	1.78	0.77
2:L:230:THR:O	2:L:231:GLU:HG3	1.85	0.76
1:A:140:ALA:HB2	3:A:448:HOH:O	1.85	0.74
2:F:123:LYS:NZ	3:F:304:HOH:O	2.10	0.74
2:F:105:GLN:O	3:F:301:HOH:O	2.03	0.74
1:E:48:LYS:HG2	1:E:49:GLY:N	2.04	0.73
2:F:75:ARG:HD2	2:F:92:SER:O	1.89	0.73
1:A:115:ARG:N	3:B:301:HOH:O	2.21	0.71
2:F:66:SER:O	3:F:303:HOH:O	2.09	0.70
1:E:48:LYS:HG2	1:E:49:GLY:H	1.56	0.69
2:D:75:ARG:HD2	2:D:92:SER:O	1.90	0.69
2:L:153:LEU:HD12	2:L:199:LEU:HD23	1.73	0.68
1:C:107:PHE:CD2	2:D:116:LEU:HD21	2.29	0.67
2:B:97:GLU:HG3	2:B:191:ASN:HD21	1.59	0.67
1:A:58:GLY:O	1:A:59:SER:OG	2.12	0.66
2:F:88:TYR:N	3:F:305:HOH:O	2.23	0.65
2:B:153:LEU:HD12	2:B:199:LEU:HD23	1.80	0.64
2:F:153:LEU:HD12	2:F:199:LEU:HD23	1.78	0.64
2:B:58:SER:HB3	2:B:87:ARG:NH2	2.13	0.64
1:A:186:GLN:NE2	3:A:301:HOH:O	2.31	0.63
1:E:45:ALA:CB	1:E:48:LYS:HE2	2.28	0.63
2:B:55:LYS:NZ	3:B:301:HOH:O	2.18	0.62
1:A:13:VAL:HG11	1:A:94:LEU:HD13	1.80	0.62
2:D:141:PRO:HD3	2:D:153:LEU:HD23	1.81	0.62
1:H:45:ALA:HB3	1:H:48:LYS:HD2	1.82	0.61
2:D:141:PRO:HD3	2:D:153:LEU:CD2	2.32	0.59
2:B:66:SER:O	2:B:67:LYS:HB2	2.03	0.58
2:L:75:ARG:HD2	2:L:92:SER:O	2.03	0.58
1:E:45:ALA:H	1:E:48:LYS:CE	2.17	0.58
2:F:4:LEU:HD21	2:F:25:LEU:HD13	1.87	0.57
1:A:59:SER:CA	1:A:82:ASN:HD21	2.05	0.57
1:A:140:ALA:HB1	1:A:141:PRO:CA	2.33	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:141:PRO:HB3	1:C:228:PRO:HA	1.87	0.56
2:D:1:GLN:N	2:D:1:GLN:OE1	2.34	0.56
2:L:141:PRO:HD3	2:L:153:LEU:CD2	2.36	0.56
1:H:114:CYS:O	2:L:55:LYS:NZ	2.33	0.55
2:L:215:GLN:NE2	2:L:224:GLU:OE2	2.33	0.55
1:E:48:LYS:NZ	1:E:49:GLY:O	2.39	0.54
1:C:201:SER:O	1:C:202:SER:OG	2.18	0.54
2:B:97:GLU:HG3	2:B:191:ASN:ND2	2.23	0.53
2:B:210:ARG:HH11	2:B:210:ARG:HG3	1.72	0.53
1:H:182:PRO:HG2	2:L:186:SER:OG	2.09	0.52
1:C:13:VAL:HG11	1:C:94:LEU:HD13	1.91	0.51
2:B:52:TYR:OH	2:B:55:LYS:HG3	2.10	0.51
1:C:14:GLN:HG3	1:E:222:VAL:HG22	1.91	0.51
2:B:141:PRO:HD3	2:B:153:LEU:CD2	2.41	0.51
1:A:107:PHE:CE2	2:B:116:LEU:HD21	2.46	0.51
2:F:56:VAL:HG11	3:F:305:HOH:O	2.11	0.51
2:B:141:PRO:HD3	2:B:153:LEU:HD23	1.93	0.50
2:F:52:TYR:OH	2:F:55:LYS:HG3	2.11	0.50
2:L:1:GLN:N	2:L:1:GLN:OE1	2.39	0.50
2:F:4:LEU:CD2	2:F:25:LEU:HD13	2.42	0.49
2:L:99:GLU:HG2	2:L:123:LYS:HZ1	1.77	0.49
2:D:129:GLN:HB2	2:D:161:TYR:CE2	2.48	0.49
1:H:174:LEU:HD21	1:H:197:VAL:HG21	1.93	0.49
2:L:168:ALA:O	2:L:214:CYS:O	2.30	0.49
1:E:45:ALA:H	1:E:48:LYS:HE2	1.76	0.49
1:H:221:LYS:NZ	1:H:223:ASP:OD1	2.45	0.49
2:B:86:GLU:OE1	2:B:88:TYR:OH	2.24	0.48
1:C:69:ASP:OD1	1:C:72:LYS:NZ	2.43	0.48
1:C:107:PHE:CE2	2:D:116:LEU:HD21	2.48	0.48
1:H:110:LYS:HD2	1:H:110:LYS:HA	1.69	0.48
2:F:187:LYS:HE3	2:F:193:TYR:CE1	2.50	0.47
2:F:215:GLN:NE2	2:F:224:GLU:OE2	2.25	0.47
1:H:229:LYS:HD2	1:H:229:LYS:HA	1.68	0.47
3:A:349:HOH:O	1:C:20:ARG:CZ	2.62	0.47
2:F:11:ALA:HB2	3:F:360:HOH:O	2.14	0.47
2:F:99:GLU:OE1	2:F:187:LYS:NZ	2.41	0.46
1:E:109:GLY:HA2	2:F:107:TRP:CZ2	2.50	0.46
2:B:48:LYS:HB3	2:B:48:LYS:HE2	1.73	0.46
1:C:52:TRP:CD1	2:D:116:LEU:HB2	2.50	0.46
2:D:39:ILE:HG12	2:D:87:ARG:HG3	1.96	0.46
1:C:116:ASP:HB2	2:D:52:TYR:CD1	2.51	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:4:LEU:HD21	2:L:25:LEU:HD13	1.98	0.45
2:F:204:GLU:CD	2:F:204:GLU:H	2.20	0.45
1:E:141:PRO:HA	1:E:152:ALA:O	2.16	0.45
1:H:107:PHE:CE2	2:L:116:LEU:HD21	2.51	0.45
2:B:2:LEU:HB2	2:B:29:HIS:CE1	2.53	0.44
2:B:66:SER:O	2:B:67:LYS:CB	2.64	0.44
2:D:153:LEU:HD12	2:D:199:LEU:HD23	1.99	0.44
1:E:48:LYS:HE3	1:E:48:LYS:HB3	1.50	0.44
1:H:13:VAL:HG11	1:H:94:LEU:HD13	1.99	0.44
1:A:141:PRO:HD2	1:A:228:PRO:HA	1.99	0.43
2:F:64:SER:HB2	3:F:400:HOH:O	2.18	0.43
2:L:2:LEU:HD13	2:L:29:HIS:CG	2.53	0.43
2:F:141:PRO:HD3	2:F:153:LEU:HD23	1.99	0.43
2:F:108:GLY:O	2:F:115:ILE:HG13	2.19	0.43
1:C:67:TYR:O	1:C:72:LYS:NZ	2.45	0.43
1:C:136:VAL:HB	1:C:222:VAL:HG11	2.01	0.43
2:F:55:LYS:N	3:F:303:HOH:O	2.25	0.43
1:C:174:LEU:HD21	1:C:197:VAL:HG21	2.01	0.43
2:F:187:LYS:HE2	2:F:191:ASN:O	2.19	0.42
1:A:214:ASN:ND2	1:A:216:LYS:HE3	2.34	0.42
2:F:99:GLU:OE2	2:F:126:VAL:N	2.40	0.42
1:E:45:ALA:H	1:E:48:LYS:NZ	2.18	0.42
2:F:106:THR:HG22	2:F:117:VAL:HG13	2.01	0.42
1:H:30:PHE:CD2	1:H:85:ASN:HA	2.54	0.42
2:L:141:PRO:HD3	2:L:153:LEU:HD21	2.01	0.42
2:D:108:GLY:O	2:D:115:ILE:HG12	2.20	0.41
2:F:39:ILE:HG21	3:F:403:HOH:O	2.20	0.41
2:L:4:LEU:CD2	2:L:25:LEU:HD13	2.50	0.41
2:B:57:ASN:OD1	2:B:59:ASP:HB3	2.20	0.41
1:E:45:ALA:HB3	1:E:48:LYS:CE	2.43	0.41
1:A:140:ALA:HB1	1:A:141:PRO:C	2.41	0.41
2:D:41:TRP:CE2	2:D:89:LEU:HB2	2.56	0.41
2:F:141:PRO:HD3	2:F:153:LEU:CD2	2.51	0.41
2:F:41:TRP:CE2	2:F:89:LEU:HB2	2.55	0.41
2:L:2:LEU:HD11	2:L:25:LEU:HD12	2.03	0.41
2:D:3:VAL:HB	2:D:26:SER:HB3	2.03	0.40
2:F:66:SER:N	3:F:303:HOH:O	2.43	0.40
2:F:80:SER:HB3	2:F:87:ARG:HD3	2.02	0.40
2:L:213:SER:HB3	2:L:225:LYS:O	2.21	0.40
1:A:115:ARG:N	3:A:310:HOH:O	2.54	0.40
2:B:72:PRO:HB2	2:B:75:ARG:HG2	2.02	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:75:ARG:HH11	1:C:75:ARG:HD2	1.77	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	202/227 (89%)	199 (98%)	2 (1%)	1 (0%)	29	11
1	C	202/227 (89%)	197 (98%)	5 (2%)	0	100	100
1	E	203/227 (89%)	199 (98%)	2 (1%)	2 (1%)	15	3
1	H	205/227 (90%)	201 (98%)	4 (2%)	0	100	100
2	B	211/216 (98%)	199 (94%)	9 (4%)	3 (1%)	11	1
2	D	211/216 (98%)	207 (98%)	4 (2%)	0	100	100
2	F	211/216 (98%)	206 (98%)	4 (2%)	1 (0%)	29	11
2	L	213/216 (99%)	206 (97%)	5 (2%)	2 (1%)	17	3
All	All	1658/1772 (94%)	1614 (97%)	35 (2%)	9 (0%)	29	11

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	140	ALA
2	B	36	SER
2	B	67	LYS
2	L	215	GLN
1	E	205	GLY
2	B	59	ASP
2	F	128	SER
2	L	30	SER
1	E	141	PRO

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	172/188 (92%)	172 (100%)	0	100	100
1	C	172/188 (92%)	172 (100%)	0	100	100
1	E	172/188 (92%)	170 (99%)	2 (1%)	71	51
1	H	174/188 (93%)	174 (100%)	0	100	100
2	B	179/182 (98%)	179 (100%)	0	100	100
2	D	179/182 (98%)	177 (99%)	2 (1%)	73	55
2	F	179/182 (98%)	179 (100%)	0	100	100
2	L	181/182 (100%)	180 (99%)	1 (1%)	86	75
All	All	1408/1480 (95%)	1403 (100%)	5 (0%)	91	84

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	177	LYS
2	D	210	ARG
1	E	115	ARG
1	E	142	SER
2	L	213	SER

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	14	GLN
1	A	186	GLN
2	B	147	GLN
1	C	120	GLN
1	E	3	GLN
1	E	207	GLN
2	F	1	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 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, 95th 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 ²)	Q<0.9
1	A	208/227 (91%)	0.01	3 (1%) 75 76	20, 26, 38, 54	0
1	C	208/227 (91%)	0.05	5 (2%) 59 58	19, 26, 51, 71	0
1	E	209/227 (92%)	0.07	4 (1%) 66 67	17, 26, 51, 67	0
1	H	211/227 (92%)	0.00	6 (2%) 53 51	17, 26, 51, 71	0
2	B	213/216 (98%)	0.41	15 (7%) 16 14	22, 34, 71, 103	0
2	D	213/216 (98%)	-0.04	6 (2%) 53 51	22, 32, 48, 59	0
2	F	213/216 (98%)	0.46	13 (6%) 21 18	20, 34, 53, 66	0
2	L	215/216 (99%)	0.18	12 (5%) 24 22	18, 31, 46, 64	0
All	All	1690/1772 (95%)	0.14	64 (3%) 40 38	17, 29, 52, 103	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	28	GLY	8.8
2	B	36	SER	8.3
2	F	67	LYS	8.0
2	B	58	SER	6.7
2	B	37	TYR	5.6
2	B	107	TRP	5.5
2	L	231	GLU	5.3
2	B	29	HIS	5.3
2	L	63	GLY	5.1
2	B	66	SER	4.5
2	F	28	GLY	4.5
1	A	205	GLY	4.2
2	F	27	SER	4.1
2	F	63	GLY	4.0
2	B	56	VAL	4.0
2	F	38	VAL	3.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	140	ALA	3.9
2	F	109	ALA	3.8
2	B	65	HIS	3.6
2	L	107	TRP	3.6
2	B	63	GLY	3.5
1	E	229	LYS	3.5
1	H	204	LEU	3.5
2	F	1	GLN	3.4
1	C	1	GLU	3.4
1	A	59	SER	3.3
2	F	71	ILE	3.2
2	B	27	SER	3.2
1	C	202	SER	3.1
1	E	141	PRO	3.0
2	B	30	SER	3.0
1	H	109	GLY	2.9
2	F	107	TRP	2.9
2	B	128	SER	2.9
1	E	142	SER	2.7
1	E	205	GLY	2.6
1	C	141	PRO	2.6
2	L	64	SER	2.6
1	C	206	THR	2.6
2	D	177	LYS	2.6
1	H	205	GLY	2.5
2	L	66	SER	2.5
2	F	58	SER	2.5
1	H	1	GLU	2.4
2	D	128	SER	2.4
2	L	214	CYS	2.4
2	D	69	ASP	2.4
2	D	63	GLY	2.4
2	B	59	ASP	2.3
1	C	149	GLY	2.3
2	F	108	GLY	2.3
1	H	150	THR	2.3
1	H	47	GLY	2.3
2	B	64	SER	2.2
2	L	153	LEU	2.2
2	F	117	VAL	2.1
2	L	115	ILE	2.1
2	L	7	SER	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	D	107	TRP	2.1
2	L	109	ALA	2.1
2	D	46	SER	2.1
2	F	66	SER	2.0
2	L	65	HIS	2.0
2	L	177	LYS	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.