



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:05 AM GMT

PDB ID : 3O11
Title : Anti-beta-amyloid antibody c706 fab in space group c2
Authors : Teplyakov, A.; Obmolova, G.; Malia, T.; Gilliland, G.L.
Deposited on : 2010-07-20
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
A user guide is available at
<http://wwpdb.org/validation/2016/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
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

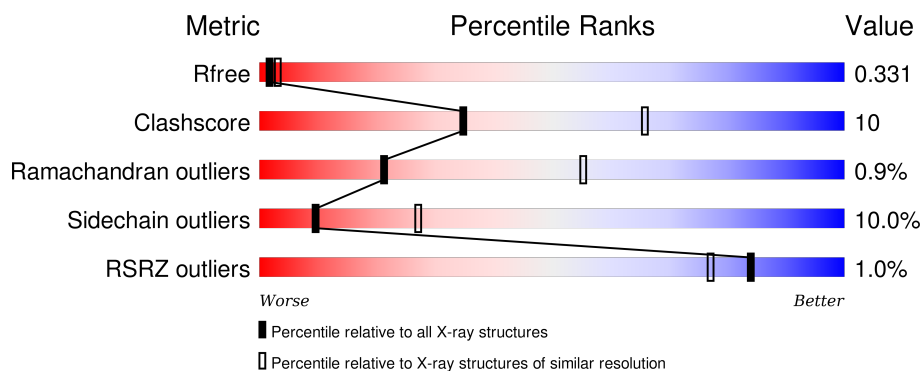
1 Overall quality at a glance

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))
R_{free}	91344	2393 (2.80-2.80)
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 ≥ 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	212	<div> <div>2%</div> <div>68%</div> <div>29%</div> <div>.</div> </div>
1	L	212	<div> <div>73%</div> <div>23%</div> <div>.</div> </div>
2	B	228	<div> <div>%</div> <div>77%</div> <div>19%</div> <div>..</div> </div>
2	H	228	<div> <div>69%</div> <div>25%</div> <div>..</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6644 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 C706 LIGHT CHAIN variable region, Ig kappa chain C region chimera.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	212	Total	C	N	O	S	0	1	0
			1614	1004	273	328	9			
1	A	212	Total	C	N	O	S	0	1	0
			1614	1004	273	328	9			

- Molecule 2 is a protein called C706 HEAVY CHAIN variable region, Ig gamma-1 chain C region chimera.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	222	Total	C	N	O	S	0	1	0
			1668	1048	288	325	7			
2	B	221	Total	C	N	O	S	0	1	0
			1652	1038	284	323	7			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	223	HIS	-	EXPRESSION TAG	UNP P01857
H	224	HIS	-	EXPRESSION TAG	UNP P01857
H	225	HIS	-	EXPRESSION TAG	UNP P01857
H	226	HIS	-	EXPRESSION TAG	UNP P01857
H	227	HIS	-	EXPRESSION TAG	UNP P01857
H	228	HIS	-	EXPRESSION TAG	UNP P01857
B	223	HIS	-	EXPRESSION TAG	UNP P01857
B	224	HIS	-	EXPRESSION TAG	UNP P01857
B	225	HIS	-	EXPRESSION TAG	UNP P01857
B	226	HIS	-	EXPRESSION TAG	UNP P01857
B	227	HIS	-	EXPRESSION TAG	UNP P01857
B	228	HIS	-	EXPRESSION TAG	UNP P01857

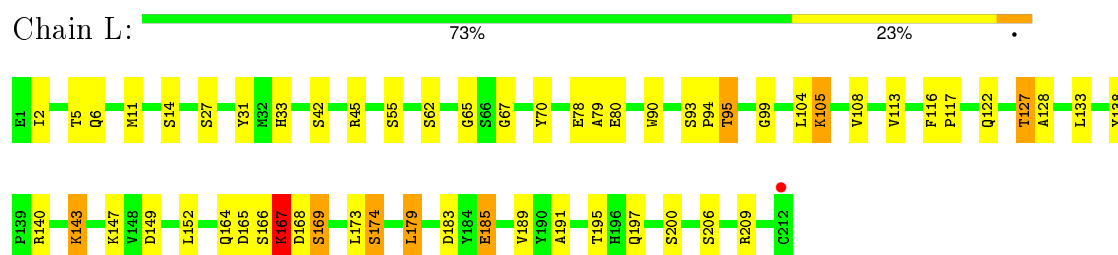
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	11	Total 11	O 11	0	0
3	H	21	Total 21	O 21	0	0
3	A	30	Total 30	O 30	0	0
3	B	34	Total 34	O 34	0	0

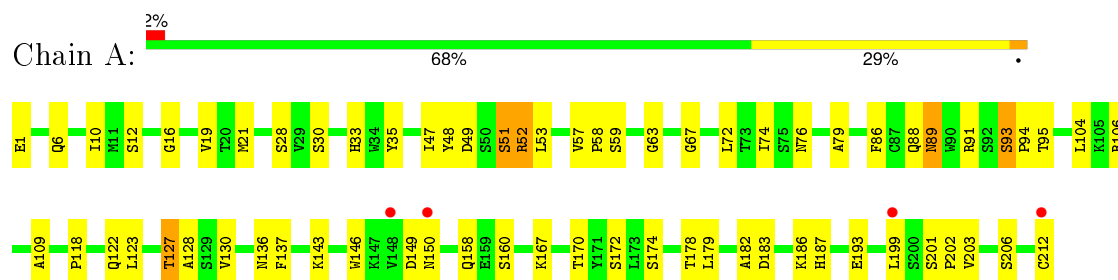
3 Residue-property plots [i](#)

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

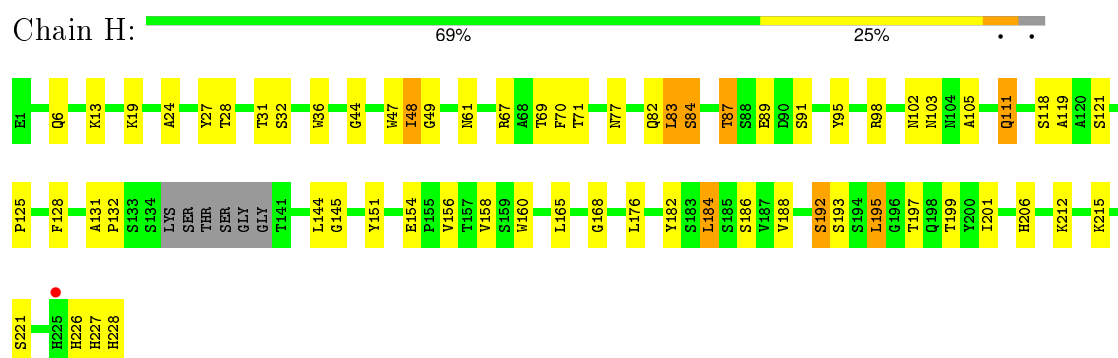
- Molecule 1: C706 LIGHT CHAIN variable region, Ig kappa chain C region chimera



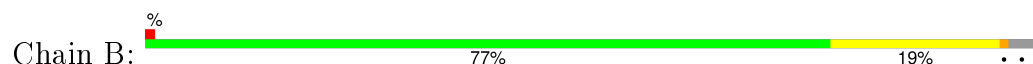
- Molecule 1: C706 LIGHT CHAIN variable region, Ig kappa chain C region chimera

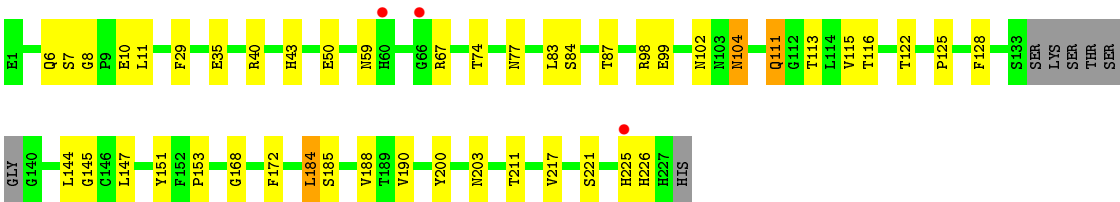


- Molecule 2: C706 HEAVY CHAIN variable region, Ig gamma-1 chain C region chimera



- Molecule 2: C706 HEAVY CHAIN variable region, Ig gamma-1 chain C region chimera





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	136.04Å 45.02Å 158.20Å 90.00° 92.69° 90.00°	Depositor
Resolution (Å)	15.00 – 2.80 42.61 – 2.80	Depositor EDS
% Data completeness (in resolution range)	83.3 (15.00-2.80) 85.3 (42.61-2.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.220 , 0.304 0.218 , 0.331	Depositor DCC
R_{free} test set	432 reflections (2.16%)	DCC
Wilson B-factor (Å ²)	73.6	Xtriage
Anisotropy	0.290	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 67.6	EDS
Estimated twinning fraction	0.000 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtriage
Outliers	2 of 20617 reflections (0.010%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6644	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.90% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PCA

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	0.49	0/1647	0.63	0/2235
1	L	0.49	0/1647	0.62	0/2235
2	B	0.48	0/1693	0.66	2/2311 (0.1%)
2	H	0.46	0/1710	0.65	0/2333
All	All	0.48	0/6697	0.64	2/9114 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	184	LEU	CA-CB-CG	6.40	130.01	115.30
2	B	226	HIS	N-CA-CB	-6.34	99.19	110.60

There are no chirality outliers.

There are no planarity outliers.

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	1614	0	1552	43	0
1	L	1614	0	1552	27	0
2	B	1652	0	1588	36	0
2	H	1668	0	1608	36	0
3	A	30	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	34	0	0	1	0
3	H	21	0	0	2	0
3	L	11	0	0	1	0
All	All	6644	0	6300	127	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 10.

All (127) 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:212:CYS:CB	2:B:225:HIS:HE1	1.17	1.53
1:A:212:CYS:HB2	2:B:225:HIS:CE1	1.46	1.50
1:A:212:CYS:CB	2:B:225:HIS:CE1	2.09	1.24
2:H:6:GLN:H	2:H:111:GLN:HE22	1.15	0.94
2:H:132:PRO:HG3	2:H:144:LEU:HB3	1.48	0.92
1:A:212:CYS:HB3	2:B:225:HIS:HE1	1.37	0.89
2:B:111:GLN:H	2:B:111:GLN:HE21	1.19	0.87
2:B:6:GLN:H	2:B:111:GLN:HE22	1.30	0.77
1:A:79:ALA:HA	1:A:104:LEU:HD13	1.71	0.73
1:L:185:GLU:HG3	1:L:209:ARG:HD2	1.71	0.73
1:A:212:CYS:SG	2:B:225:HIS:HE1	2.13	0.71
1:A:212:CYS:HB3	2:B:225:HIS:CE1	2.14	0.71
1:A:212:CYS:SG	2:B:225:HIS:CE1	2.84	0.71
2:H:156:VAL:HG12	2:H:206:HIS:HB2	1.77	0.66
1:L:127:THR:HG22	1:L:128:ALA:H	1.61	0.66
1:A:89:ASN:ND2	1:A:91:ARG:H	1.94	0.66
2:B:29:PHE:HB2	2:B:77:ASN:OD1	1.97	0.64
2:H:49:GLY:HA3	2:H:70:PHE:CE2	2.32	0.64
1:L:80:GLU:HG2	3:L:213:HOH:O	1.97	0.64
1:L:31:TYR:OH	1:L:90:TRP:HB2	1.99	0.62
1:A:212:CYS:HB2	2:B:225:HIS:NE2	2.09	0.61
1:L:128:ALA:HB3	1:L:179:LEU:O	2.02	0.59
2:B:8:GLY:O	2:B:113:THR:HG23	2.03	0.59
2:H:67:ARG:HA	2:H:84:SER:HB2	1.86	0.58
1:L:14:SER:HB3	1:L:105:LYS:HD2	1.86	0.57
1:A:118:PRO:HD3	1:A:130:VAL:HG22	1.85	0.57
1:L:65:GLY:HA3	1:L:70:TYR:HA	1.86	0.56
2:H:111:GLN:HE21	2:H:111:GLN:H	1.52	0.56
2:B:67:ARG:NH1	2:B:84:SER:O	2.40	0.55
1:A:21:MET:HG2	1:A:72:LEU:HB3	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:ASN:HD22	1:A:91:ARG:H	1.53	0.55
1:L:2:ILE:HG12	1:L:27:SER:HB3	1.88	0.55
2:B:40:ARG:HH21	2:B:43:HIS:CE1	2.24	0.55
2:B:111:GLN:HG3	3:B:256:HOH:O	2.07	0.54
1:A:49:ASP:HB2	1:A:52:ARG:HG3	1.89	0.54
1:L:147:LYS:HB2	1:L:191:ALA:HB3	1.90	0.54
2:H:160:TRP:HB3	2:H:165:LEU:HD23	1.88	0.54
2:H:13:LYS:HG2	2:H:119:ALA:HA	1.90	0.53
1:L:93:SER:HA	1:L:94:PRO:C	2.28	0.53
2:B:98:ARG:HD2	2:B:99:GLU:O	2.09	0.53
1:A:93:SER:HA	1:A:94:PRO:C	2.27	0.52
2:B:125:PRO:HD2	2:B:211:THR:HG21	1.90	0.52
1:L:167:LYS:HE3	1:L:167:LYS:HA	1.91	0.52
2:H:125:PRO:HB3	2:H:151:TYR:HB3	1.91	0.52
2:B:6:GLN:H	2:B:111:GLN:NE2	2.04	0.52
1:L:33:HIS:CE1	2:H:105:ALA:HB2	2.44	0.52
1:A:172:SER:O	2:B:172:PHE:HE1	1.94	0.51
1:A:28:SER:HB2	3:A:216:HOH:O	2.09	0.51
2:B:111:GLN:N	2:B:111:GLN:HE21	1.99	0.51
1:A:47:ILE:HG12	1:A:53:LEU:HD23	1.93	0.50
1:L:143:LYS:HB3	1:L:195:THR:HB	1.93	0.50
2:B:168:GLY:O	2:B:188:VAL:HA	2.12	0.50
2:B:50:GLU:OE2	2:B:59:ASN:ND2	2.43	0.50
2:B:11:LEU:HD23	2:B:122:THR:OG1	2.11	0.50
1:A:53:LEU:HD22	1:A:57:VAL:HG11	1.94	0.49
2:H:47:TRP:CZ2	2:H:49:GLY:HA2	2.47	0.49
1:A:1:PCA:HB2	1:A:93:SER:OG	2.12	0.49
2:H:132:PRO:CG	2:H:144:LEU:HB3	2.32	0.49
2:H:227:HIS:HB3	2:H:228:HIS:ND1	2.26	0.49
2:B:147:LEU:HD12	2:B:185:SER:HB3	1.93	0.48
2:H:168:GLY:O	2:H:188:VAL:HA	2.14	0.48
2:B:35:GLU:OE2	2:B:50:GLU:CB	2.61	0.48
1:A:199:LEU:HD13	1:A:203:VAL:HG23	1.96	0.48
2:H:36:TRP:HB3	2:H:48:ILE:HD12	1.95	0.48
1:A:106:ARG:NH2	1:A:109:ALA:HB2	2.29	0.47
2:H:184:LEU:HD12	2:H:184:LEU:C	2.35	0.47
2:H:226:HIS:O	2:H:227:HIS:HD2	1.97	0.47
2:H:158:VAL:HG11	2:H:186:SER:HB3	1.96	0.47
1:A:149:ASP:OD2	1:A:187:HIS:HD2	1.98	0.47
1:A:137:PHE:O	1:A:137:PHE:CD1	2.68	0.47
2:H:36:TRP:O	2:H:48:ILE:HG13	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:24:ALA:HB1	2:H:27:TYR:CE2	2.49	0.46
2:H:154:GLU:HB2	3:H:244:HOH:O	2.14	0.46
1:L:168:ASP:O	1:L:169:SER:HB2	2.15	0.46
2:H:158:VAL:HG11	2:H:186:SER:CB	2.45	0.46
1:L:165:ASP:O	1:L:169:SER:HA	2.16	0.46
1:L:149:ASP:HA	1:L:189:VAL:HB	1.97	0.46
2:H:192:SER:HA	2:H:195:LEU:HG	1.98	0.45
1:A:193:GLU:HA	1:A:203:VAL:O	2.16	0.45
1:A:136:ASN:HB3	1:A:170:THR:HG21	1.97	0.45
1:A:16:GLY:HA2	1:A:76:ASN:HB2	1.97	0.45
2:B:125:PRO:HB3	2:B:151:TYR:HB3	1.98	0.45
1:A:122:GLN:HG3	2:B:128:PHE:CE2	2.51	0.45
1:A:149:ASP:OD2	1:A:187:HIS:CD2	2.69	0.45
2:H:83:LEU:HA	2:H:83:LEU:HD23	1.81	0.45
2:B:11:LEU:HD22	2:B:153:PRO:HD3	1.99	0.44
1:L:79:ALA:HA	1:L:104:LEU:HD13	1.99	0.44
2:H:131:ALA:HA	2:H:132:PRO:HD3	1.90	0.44
1:A:33:HIS:CE1	1:A:49:ASP:H	2.34	0.44
2:B:50:GLU:HG2	2:B:59:ASN:HB2	1.97	0.44
2:B:104:ASN:C	2:B:104:ASN:HD22	2.20	0.44
2:H:87:THR:OG1	2:H:89:GLU:HG2	2.18	0.44
1:L:173:LEU:HD23	1:L:174:SER:N	2.32	0.44
2:B:35:GLU:OE2	2:B:50:GLU:HB3	2.17	0.44
1:L:95:THR:HG21	2:H:47:TRP:HB3	2.01	0.43
1:A:118:PRO:HB2	1:A:123:LEU:HD21	1.99	0.43
1:L:104:LEU:O	1:L:164:GLN:NE2	2.51	0.43
2:H:47:TRP:O	2:H:61:ASN:ND2	2.39	0.43
1:A:33:HIS:HD1	1:A:48:TYR:HA	1.84	0.43
1:A:182:ALA:O	1:A:186:LYS:HG3	2.18	0.43
1:A:35:TYR:HE1	1:A:88:GLN:CB	2.32	0.43
2:B:35:GLU:OE2	2:B:50:GLU:HB2	2.19	0.42
1:L:6:GLN:OE1	1:L:99:GLY:HA2	2.19	0.42
2:H:44:GLY:HA2	3:H:231:HOH:O	2.19	0.42
1:A:33:HIS:NE2	2:B:104:ASN:ND2	2.67	0.42
1:L:42:SER:HA	2:H:95:TYR:CE1	2.55	0.42
1:A:19:VAL:HB	1:A:74:ILE:HD12	2.01	0.42
1:A:6:GLN:OE1	1:A:86:PHE:HA	2.19	0.42
1:L:116:PHE:HA	1:L:117:PRO:HD3	1.87	0.42
1:L:122:GLN:HG3	2:H:128:PHE:CE2	2.54	0.42
1:A:127:THR:HG22	1:A:128:ALA:H	1.84	0.42
2:B:145:GLY:O	2:B:217:VAL:HG21	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:176:LEU:HD13	2:H:182:TYR:CZ	2.55	0.41
2:H:19:LYS:HG3	2:H:82:GLN:HB2	2.01	0.41
2:B:190:VAL:HG11	2:B:200:TYR:CE1	2.56	0.41
1:L:113:VAL:HA	1:L:133:LEU:O	2.20	0.41
1:A:51:SER:HB3	1:A:63:GLY:O	2.21	0.41
1:L:185:GLU:HA	1:L:209:ARG:NE	2.36	0.41
2:H:201:ILE:HA	2:H:215:LYS:O	2.21	0.41
2:H:145:GLY:HA2	2:H:160:TRP:CH2	2.55	0.41
2:B:10:GLU:HB2	2:B:115:VAL:HG22	2.03	0.41
2:H:28:THR:HB	2:H:31:THR:OG1	2.20	0.41
1:L:108:VAL:HG13	1:L:138:TYR:O	2.21	0.41
1:A:53:LEU:HD22	1:A:57:VAL:CG1	2.51	0.40
1:A:57:VAL:HA	1:A:58:PRO:HD3	1.97	0.40
1:A:193:GLU:OE2	1:A:202:PRO:HB3	2.22	0.40
1:A:130:VAL:HG12	1:A:146:TRP:CH2	2.57	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	211/212 (100%)	199 (94%)	11 (5%)	1 (0%)	34	69
1	L	211/212 (100%)	187 (89%)	21 (10%)	3 (1%)	14	42
2	B	218/228 (96%)	197 (90%)	20 (9%)	1 (0%)	34	69
2	H	219/228 (96%)	197 (90%)	19 (9%)	3 (1%)	14	42
All	All	859/880 (98%)	780 (91%)	71 (8%)	8 (1%)	21	55

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	67	GLY
1	L	167	LYS
2	H	91	SER
2	H	195	LEU
2	B	221	SER
1	L	169	SER
2	H	221	SER
1	A	67	GLY

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	183/183 (100%)	162 (88%)	21 (12%)	7	21
1	L	183/183 (100%)	161 (88%)	22 (12%)	6	19
2	B	183/189 (97%)	172 (94%)	11 (6%)	24	56
2	H	186/189 (98%)	166 (89%)	20 (11%)	8	23
All	All	735/744 (99%)	661 (90%)	74 (10%)	9	27

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	5	THR
1	L	11	MET
1	L	45	ARG
1	L	55	SER
1	L	62	SER
1	L	78	GLU
1	L	95	THR
1	L	105	LYS
1	L	127	THR
1	L	140	ARG
1	L	143	LYS
1	L	152	LEU
1	L	166[A]	SER
1	L	166[B]	SER

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Mol	Chain	Res	Type
1	L	167	LYS
1	L	174	SER
1	L	179	LEU
1	L	183	ASP
1	L	185	GLU
1	L	197	GLN
1	L	200	SER
1	L	206	SER
2	H	32	SER
2	H	48	ILE
2	H	69	THR
2	H	71	THR
2	H	77	ASN
2	H	83	LEU
2	H	84	SER
2	H	87	THR
2	H	98	ARG
2	H	102	ASN
2	H	103	ASN
2	H	111	GLN
2	H	118	SER
2	H	121	SER
2	H	184	LEU
2	H	192	SER
2	H	193	SER
2	H	197	THR
2	H	199	THR
2	H	212	LYS
1	A	10	ILE
1	A	12	SER
1	A	30	SER
1	A	51	SER
1	A	52	ARG
1	A	59	SER
1	A	89	ASN
1	A	93	SER
1	A	95	THR
1	A	127	THR
1	A	143	LYS
1	A	150	ASN
1	A	158	GLN
1	A	160	SER

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Mol	Chain	Res	Type
1	A	167	LYS
1	A	174	SER
1	A	178	THR
1	A	179	LEU
1	A	183	ASP
1	A	201	SER
1	A	206	SER
2	B	7	SER
2	B	74	THR
2	B	83	LEU
2	B	87	THR
2	B	102	ASN
2	B	104	ASN
2	B	111	GLN
2	B	116	THR
2	B	144	LEU
2	B	184	LEU
2	B	203	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	37	GLN
1	L	89	ASN
1	L	136	ASN
1	L	158	GLN
1	L	197	GLN
2	H	3	GLN
2	H	111	GLN
2	H	170	HIS
2	H	225	HIS
2	H	227	HIS
1	A	76	ASN
1	A	89	ASN
1	A	145	GLN
1	A	187	HIS
2	B	3	GLN
2	B	104	ASN
2	B	111	GLN
2	B	225	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	PCA	A	1	1	7,8,9	1.62	1 (14%)	9,10,12	2.17	4 (44%)
2	PCA	B	1	2	7,8,9	1.48	1 (14%)	9,10,12	2.09	4 (44%)
2	PCA	H	1	2	7,8,9	1.57	1 (14%)	9,10,12	2.05	4 (44%)
1	PCA	L	1	1	7,8,9	1.65	1 (14%)	9,10,12	2.22	5 (55%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	PCA	A	1	1	-	0/0/11/13	0/1/1/1
2	PCA	B	1	2	-	0/0/11/13	0/1/1/1
2	PCA	H	1	2	-	0/0/11/13	0/1/1/1
1	PCA	L	1	1	-	0/0/11/13	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	PCA	CD-N	3.73	1.46	1.33
1	A	1	PCA	CD-N	3.75	1.46	1.33
2	H	1	PCA	CD-N	3.99	1.47	1.33
1	L	1	PCA	CD-N	4.19	1.47	1.33

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	PCA	CA-N-CD	-4.25	99.58	113.81
1	A	1	PCA	CA-N-CD	-4.23	99.62	113.81
1	L	1	PCA	CB-CA-C	-3.70	107.71	112.76
2	H	1	PCA	CA-N-CD	-3.11	103.39	113.81
2	H	1	PCA	CB-CA-C	-3.09	108.54	112.76
1	L	1	PCA	CA-N-CD	-2.92	104.01	113.81
1	A	1	PCA	O-C-CA	-2.73	118.23	125.44
1	L	1	PCA	OE-CD-CG	-2.63	120.95	126.81
2	H	1	PCA	OE-CD-CG	-2.36	121.53	126.81
1	L	1	PCA	O-C-CA	-2.25	119.51	125.44
2	B	1	PCA	CB-CA-C	-2.17	109.80	112.76
2	H	1	PCA	O-C-CA	-2.16	119.73	125.44
1	A	1	PCA	CB-CA-C	-2.16	109.82	112.76
2	B	1	PCA	OE-CD-CG	-2.01	122.32	126.81
1	A	1	PCA	CB-CA-N	2.10	109.32	103.20
2	B	1	PCA	CB-CA-N	2.10	109.34	103.20
1	L	1	PCA	CB-CA-N	2.14	109.46	103.20

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1	PCA	1	0

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

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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	211/212 (99%)	0.04	4 (1%) 70 59	45, 69, 92, 131	0
1	L	211/212 (99%)	0.13	1 (0%) 91 88	45, 71, 97, 139	0
2	B	220/228 (96%)	-0.06	3 (1%) 78 69	34, 66, 97, 127	0
2	H	221/228 (96%)	-0.00	1 (0%) 91 88	45, 70, 100, 125	0
All	All	863/880 (98%)	0.03	9 (1%) 84 77	34, 69, 97, 139	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	225	HIS	3.6
2	H	225	HIS	3.5
1	A	212	CYS	3.4
1	L	212	CYS	3.2
2	B	60	HIS	2.3
1	A	148	VAL	2.2
1	A	150	ASN	2.1
1	A	199	LEU	2.1
2	B	66	GLY	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	PCA	L	1	8/9	0.93	0.16	-	75,79,82,82	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	PCA	B	1	8/9	0.94	0.17	-	70,73,77,82	0
1	PCA	A	1	8/9	0.76	0.36	-	80,81,82,83	0
2	PCA	H	1	8/9	0.91	0.23	-	78,87,89,98	0

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.