



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:28 AM GMT

PDB ID : 3AUV
Title : Predicting Amino Acid Preferences in the Complementarity Determining Regions of an Antibody-Antigen Recognition Interface
Authors : Yu, C.M.; Peng, H.P.; Chen, I.C.; Lee, Y.C.; Chen, J.B.; Tsai, K.C.; Chen, C.T.; Chang, J.Y.; Yang, E.W.; Hsu, P.C.; Jian, J.W.; Hsu, H.J.; Chang, H.J.; Hsu, W.L.; Huang, K.F.; Ma, A.C.; Yang, A.S.
Deposited on : 2011-02-16
Resolution : 2.40 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

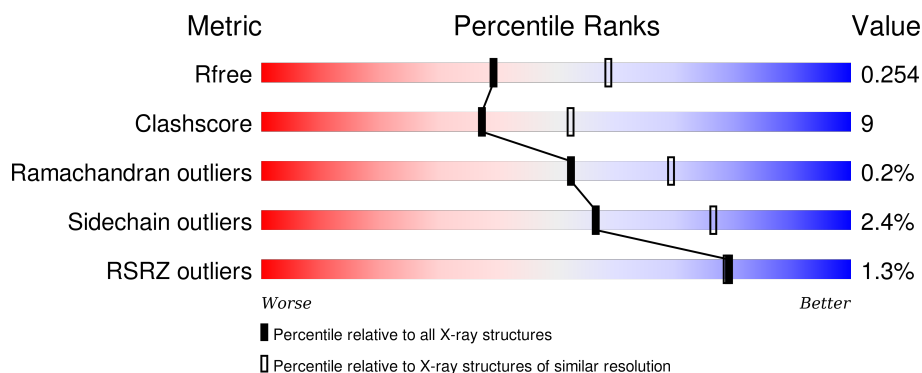
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	276	<div> <div>2%</div> <div>68% 13% 18%</div> </div>
1	B	276	<div> <div>65% 15% 18%</div> </div>
1	C	276	<div> <div>2%</div> <div>67% 13% 18%</div> </div>
1	D	276	<div> <div>66% 15% 18%</div> </div>
1	E	276	<div> <div>2%</div> <div>67% 14% 18%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	276	<div><div><div>%</div><div><div></div></div><div>67%</div><div>13%</div><div>19%</div></div></div>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 10863 atoms, of which 0 are hydrogens and 0 are deuteriums.

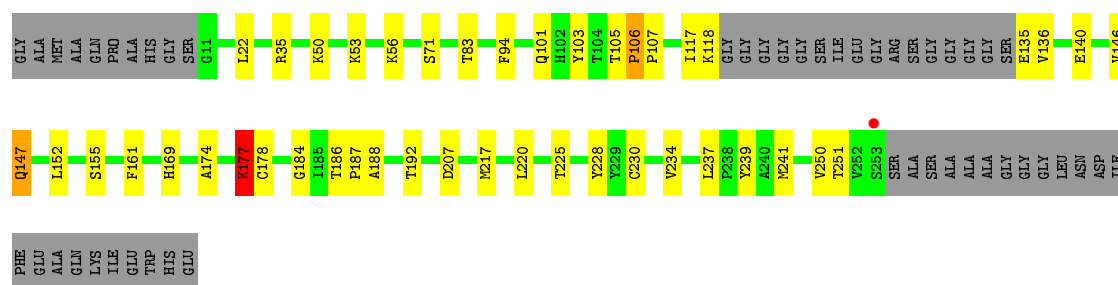
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called sc-dsFv derived from the G6-Fab.

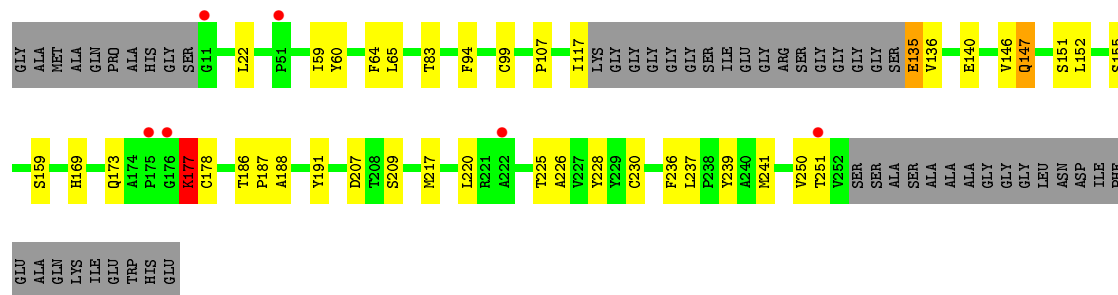
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	225	Total	C	N	O	S	0	0	0
			1721	1098	280	334	9			
1	B	225	Total	C	N	O	S	0	0	0
			1719	1097	280	333	9			
1	C	225	Total	C	N	O	S	0	0	0
			1719	1097	280	333	9			
1	D	227	Total	C	N	O	S	0	0	0
			1734	1106	283	336	9			
1	E	225	Total	C	N	O	S	0	0	0
			1719	1097	280	333	9			
1	F	224	Total	C	N	O	S	0	0	0
			1710	1092	279	330	9			

- Molecule 2 is water.

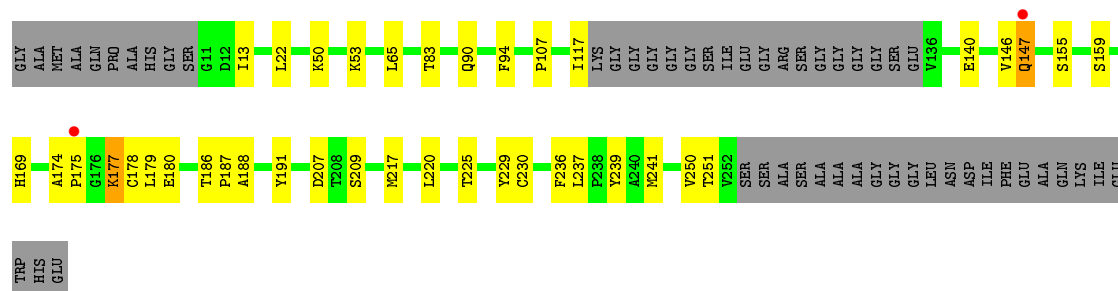
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	92	Total	O	0	0
			92	92		
2	B	99	Total	O	0	0
			99	99		
2	C	93	Total	O	0	0
			93	93		
2	D	94	Total	O	0	0
			94	94		
2	E	86	Total	O	0	0
			86	86		
2	F	77	Total	O	0	0
			77	77		



- Molecule 1: sc-dsFv derived from the G6-Fab



- Molecule 1: sc-dsFv derived from the G6-Fab



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	136.06Å 136.06Å 169.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.32 – 2.40 29.46 – 2.40	Depositor EDS
% Data completeness (in resolution range)	89.9 (29.32-2.40) 99.1 (29.46-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.70 (at 2.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_648)	Depositor
R, R_{free}	0.216 , 0.251 0.220 , 0.254	Depositor DCC
R_{free} test set	3536 reflections (5.06%)	DCC
Wilson B-factor (Å ²)	28.1	Xtriage
Anisotropy	0.786	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 38.9	EDS
Estimated twinning fraction	0.000 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.37$	Xtriage
Outliers	27 of 70368 reflections (0.038%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10863	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.15 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.3770e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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	0.90	0/1767	0.86	1/2408 (0.0%)
1	B	0.86	0/1765	0.83	1/2405 (0.0%)
1	C	0.90	0/1765	0.86	0/2405
1	D	0.86	0/1780	0.86	1/2424 (0.0%)
1	E	0.92	1/1765 (0.1%)	0.91	4/2405 (0.2%)
1	F	0.94	0/1756	0.87	0/2393
All	All	0.90	1/10598 (0.0%)	0.87	7/14440 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	99	CYS	CB-SG	-5.36	1.73	1.81

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	177	LYS	CD-CE-NZ	8.49	131.23	111.70
1	E	177	LYS	CA-CB-CG	6.11	126.84	113.40
1	E	177	LYS	CB-CG-CD	-6.04	95.88	111.60
1	A	72	ARG	NE-CZ-NH1	-5.77	117.42	120.30
1	B	177	LYS	CD-CE-NZ	5.60	124.57	111.70
1	E	177	LYS	N-CA-CB	5.31	120.16	110.60
1	E	177	LYS	CB-CA-C	-5.25	99.89	110.40

There are no chirality outliers.

There are no planarity outliers.

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	A	1721	0	1645	27	0
1	B	1719	0	1643	34	0
1	C	1719	0	1643	30	0
1	D	1734	0	1661	35	0
1	E	1719	0	1643	34	0
1	F	1710	0	1637	36	0
2	A	92	0	0	1	0
2	B	99	0	0	3	0
2	C	93	0	0	1	0
2	D	94	0	0	5	0
2	E	86	0	0	2	0
2	F	77	0	0	4	0
All	All	10863	0	9872	187	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 9.

All (187) 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:D:440:HOH:O	1:E:135:GLU:HB2	1.71	0.91
1:B:94:PHE:HB3	1:B:117:ILE:HD12	1.55	0.88
1:F:237:LEU:HD23	1:F:239:TYR:CE2	2.09	0.88
1:B:237:LEU:HD23	1:B:239:TYR:CE2	2.09	0.87
1:E:225:THR:HG23	1:E:251:THR:HA	1.64	0.79
1:C:225:THR:HG23	1:C:251:THR:HA	1.65	0.79
1:F:180:GLU:OE2	2:F:459:HOH:O	2.00	0.78
1:F:225:THR:HG23	1:F:251:THR:HA	1.67	0.77
1:D:225:THR:HG23	1:D:251:THR:HA	1.67	0.76
1:F:94:PHE:HB3	1:F:117:ILE:HD12	1.67	0.76
1:B:237:LEU:HD23	1:B:239:TYR:HE2	1.49	0.75
1:A:237:LEU:HD23	1:A:239:TYR:CE2	2.23	0.74
1:C:237:LEU:HD23	1:C:239:TYR:CE2	2.23	0.73
1:E:94:PHE:HB3	1:E:117:ILE:HD12	1.70	0.73
1:A:94:PHE:HB3	1:A:117:ILE:HD12	1.71	0.73
1:E:237:LEU:HD23	1:E:239:TYR:CE2	2.24	0.72
1:C:94:PHE:HB3	1:C:117:ILE:HD12	1.70	0.72
1:B:153:ARG:O	2:B:336:HOH:O	2.08	0.72
1:D:71:SER:O	2:D:466:HOH:O	2.09	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:135:GLU:HA	1:D:135:GLU:OE1	1.91	0.70
1:B:225:THR:HG23	1:B:251:THR:HA	1.72	0.70
1:D:94:PHE:HB3	1:D:117:ILE:HD12	1.73	0.70
1:F:237:LEU:HD23	1:F:239:TYR:HE2	1.53	0.69
1:C:237:LEU:HD12	1:E:207:ASP:HB2	1.76	0.68
1:E:155:SER:OG	2:E:285:HOH:O	1.91	0.68
1:D:237:LEU:HD23	1:D:239:TYR:CE2	2.28	0.68
1:F:237:LEU:CD2	1:F:239:TYR:HE2	2.07	0.67
1:C:153:ARG:O	2:C:307:HOH:O	2.13	0.66
1:A:225:THR:HG23	1:A:251:THR:HA	1.77	0.66
1:E:146:VAL:HG23	1:E:147:GLN:N	2.10	0.66
1:D:237:LEU:HD12	1:F:207:ASP:HB2	1.78	0.64
1:C:217:MET:HE2	1:C:220:LEU:HD21	1.81	0.63
1:B:237:LEU:CD2	1:B:239:TYR:HE2	2.13	0.61
1:C:59:ILE:HD13	1:C:65:LEU:HD23	1.84	0.60
1:C:146:VAL:HG23	1:C:147:GLN:N	2.17	0.59
1:E:209:SER:HA	1:F:159:SER:HB3	1.85	0.59
1:F:177:LYS:HD2	1:F:178:CYS:H	1.68	0.59
1:F:174:ALA:HB3	1:F:177:LYS:CB	2.33	0.59
1:F:13:ILE:O	2:F:281:HOH:O	2.17	0.58
1:D:118:LYS:HA	2:D:474:HOH:O	2.04	0.58
1:D:187:PRO:O	1:D:188:ALA:HB3	2.03	0.57
1:A:177:LYS:HD2	1:A:178:CYS:H	1.68	0.57
1:F:175:PRO:HG2	2:F:441:HOH:O	2.05	0.57
1:C:174:ALA:HB3	1:C:177:LYS:CB	2.35	0.57
1:B:11:GLY:N	2:B:530:HOH:O	2.37	0.57
1:D:217:MET:HE2	1:D:220:LEU:HD21	1.85	0.56
1:F:237:LEU:HD23	1:F:239:TYR:CD2	2.40	0.56
1:E:169:HIS:O	1:E:230:CYS:HA	2.04	0.56
1:C:187:PRO:O	1:C:188:ALA:HB3	2.04	0.56
1:F:90:GLN:OE1	1:F:90:GLN:HA	2.05	0.56
1:B:107:PRO:HG3	1:B:241:MET:CE	2.35	0.56
1:D:251:THR:OG1	2:D:389:HOH:O	2.18	0.55
1:B:146:VAL:HG23	1:B:147:GLN:N	2.21	0.55
1:B:187:PRO:O	1:B:188:ALA:HB3	2.06	0.55
1:B:107:PRO:HG3	1:B:241:MET:HE1	1.89	0.55
1:B:237:LEU:HD12	1:D:207:ASP:HB2	1.88	0.55
1:E:159:SER:HB3	1:F:209:SER:HA	1.89	0.55
1:F:175:PRO:O	2:F:441:HOH:O	2.18	0.54
1:A:60:TYR:O	1:A:64:PHE:HB2	2.07	0.54
1:A:217:MET:HE2	1:A:220:LEU:HD21	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:237:LEU:CD2	1:F:239:TYR:CE2	2.84	0.54
1:F:146:VAL:HG23	1:F:147:GLN:N	2.22	0.54
1:D:174:ALA:HB3	1:D:177:LYS:CB	2.38	0.54
1:E:187:PRO:O	1:E:188:ALA:HB3	2.07	0.54
1:B:207:ASP:HB2	1:F:237:LEU:HD12	1.88	0.53
1:F:174:ALA:HB3	1:F:177:LYS:HB2	1.90	0.53
1:C:177:LYS:HD2	1:C:178:CYS:H	1.74	0.53
1:D:107:PRO:HG3	1:D:241:MET:CE	2.38	0.53
1:A:187:PRO:O	1:A:188:ALA:HB3	2.08	0.53
1:F:217:MET:HE2	1:F:220:LEU:HD21	1.90	0.53
1:B:169:HIS:O	1:B:230:CYS:HA	2.09	0.53
1:E:217:MET:HE2	1:E:220:LEU:HD21	1.89	0.53
1:E:22:LEU:C	1:E:22:LEU:HD12	2.29	0.53
1:D:146:VAL:HG23	1:D:147:GLN:N	2.23	0.52
1:A:174:ALA:HB3	1:A:177:LYS:CB	2.39	0.52
1:A:169:HIS:O	1:A:230:CYS:HA	2.10	0.52
1:B:37:SER:OG	2:B:281:HOH:O	2.17	0.52
1:A:207:ASP:HB2	1:E:237:LEU:HD12	1.91	0.52
1:F:140:GLU:HA	1:F:155:SER:O	2.10	0.52
1:A:146:VAL:HG23	1:A:147:GLN:N	2.24	0.52
1:C:174:ALA:HB3	1:C:177:LYS:HB2	1.91	0.52
1:F:22:LEU:C	1:F:22:LEU:HD12	2.31	0.52
1:F:187:PRO:O	1:F:188:ALA:HB3	2.08	0.52
1:B:59:ILE:HD13	1:B:65:LEU:HD23	1.92	0.51
1:F:177:LYS:HD2	1:F:178:CYS:N	2.26	0.51
1:E:237:LEU:HD23	1:E:239:TYR:CD2	2.45	0.51
1:A:174:ALA:HB3	1:A:177:LYS:HB2	1.93	0.51
1:A:191:TYR:OH	1:E:186:THR:HG21	2.10	0.51
1:D:174:ALA:HB3	1:D:177:LYS:HB2	1.93	0.51
1:B:217:MET:HE2	1:B:220:LEU:HD21	1.92	0.51
1:D:169:HIS:O	1:D:230:CYS:HA	2.11	0.51
1:B:146:VAL:HG11	1:B:152:LEU:HD22	1.93	0.50
1:F:225:THR:HA	1:F:250:VAL:O	2.12	0.50
1:C:140:GLU:HA	1:C:155:SER:O	2.11	0.50
1:F:107:PRO:HG3	1:F:241:MET:CE	2.41	0.50
1:F:169:HIS:O	1:F:230:CYS:HA	2.12	0.50
1:A:140:GLU:HA	1:A:155:SER:O	2.12	0.49
1:E:140:GLU:HA	1:E:155:SER:O	2.12	0.49
1:E:237:LEU:CD2	1:E:239:TYR:CE2	2.95	0.49
1:B:140:GLU:OE2	1:B:228:TYR:O	2.30	0.49
1:E:225:THR:HA	1:E:250:VAL:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:169:HIS:O	1:C:230:CYS:HA	2.12	0.49
1:C:107:PRO:HG3	1:C:241:MET:CE	2.42	0.49
1:E:107:PRO:HG3	1:E:241:MET:CE	2.43	0.49
1:D:237:LEU:HD23	1:D:239:TYR:HE2	1.73	0.49
1:B:60:TYR:O	1:B:64:PHE:HB2	2.13	0.49
1:D:105:THR:OG1	1:D:106:PRO:HA	2.13	0.48
1:A:135:GLU:HG2	1:A:136:VAL:N	2.29	0.48
1:B:174:ALA:HB3	1:B:177:LYS:CB	2.43	0.48
1:D:22:LEU:C	1:D:22:LEU:HD12	2.35	0.48
1:D:140:GLU:OE2	1:D:228:TYR:O	2.33	0.47
1:E:237:LEU:CD2	1:E:239:TYR:HE2	2.27	0.47
1:B:65:LEU:HD23	1:B:65:LEU:HA	1.63	0.47
1:A:177:LYS:HD3	1:A:177:LYS:HA	1.77	0.47
1:D:56:LYS:HG2	2:D:346:HOH:O	2.14	0.47
1:F:50:LYS:HB2	1:F:53:LYS:HG3	1.96	0.47
1:E:186:THR:HB	1:E:187:PRO:HD2	1.97	0.46
1:D:107:PRO:HG3	1:D:241:MET:HE1	1.98	0.46
1:C:179:LEU:HA	1:C:179:LEU:HD23	1.72	0.46
1:D:136:VAL:HG13	1:D:161:PHE:CD1	2.50	0.46
1:C:105:THR:OG1	1:C:106:PRO:HA	2.15	0.46
1:B:140:GLU:HA	1:B:155:SER:O	2.15	0.46
1:B:237:LEU:CD2	1:B:239:TYR:CE2	2.89	0.46
1:C:146:VAL:CG2	1:C:147:GLN:N	2.79	0.45
1:A:175:PRO:HG2	2:A:308:HOH:O	2.15	0.45
1:E:65:LEU:HA	1:E:65:LEU:HD23	1.57	0.45
1:A:107:PRO:HG3	1:A:241:MET:CE	2.45	0.45
1:D:101:GLN:OE1	1:D:103:TYR:N	2.45	0.45
1:A:135:GLU:HG2	1:A:136:VAL:H	1.81	0.45
1:F:65:LEU:HA	1:F:65:LEU:HD23	1.57	0.45
1:C:22:LEU:C	1:C:22:LEU:HD12	2.37	0.45
1:F:140:GLU:OE1	1:F:229:TYR:HA	2.16	0.45
1:E:225:THR:O	1:E:226:ALA:HB2	2.18	0.44
1:E:140:GLU:OE2	1:E:228:TYR:O	2.35	0.44
1:A:50:LYS:HB2	1:A:53:LYS:HG3	2.00	0.44
1:A:65:LEU:HD23	1:A:65:LEU:HA	1.51	0.44
1:D:146:VAL:HG11	1:D:152:LEU:HD22	2.00	0.44
1:E:236:PHE:O	1:E:239:TYR:N	2.50	0.44
1:B:22:LEU:C	1:B:22:LEU:HD12	2.38	0.44
1:B:179:LEU:HA	1:B:179:LEU:HD23	1.77	0.44
1:C:177:LYS:HD2	1:C:178:CYS:N	2.33	0.44
1:A:177:LYS:HD2	1:A:178:CYS:N	2.31	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:186:THR:HG21	1:F:191:TYR:OH	2.17	0.43
1:E:186:THR:HB	1:E:187:PRO:CD	2.48	0.43
1:F:107:PRO:HG3	1:F:241:MET:HE1	2.01	0.43
1:F:179:LEU:HA	1:F:179:LEU:HD23	1.75	0.43
1:D:50:LYS:HB2	1:D:53:LYS:HG3	2.01	0.43
1:B:41:SER:OG	1:B:42:THR:N	2.50	0.43
1:C:225:THR:HA	1:C:250:VAL:O	2.18	0.43
1:E:173:GLN:O	1:E:226:ALA:HB1	2.17	0.43
1:A:225:THR:HA	1:A:250:VAL:O	2.19	0.43
1:A:59:ILE:HD13	1:A:65:LEU:HD23	2.01	0.43
1:D:140:GLU:HA	1:D:155:SER:O	2.19	0.43
1:D:225:THR:HA	1:D:250:VAL:O	2.18	0.43
1:C:65:LEU:HD23	1:C:65:LEU:HA	1.59	0.43
1:B:225:THR:HA	1:B:250:VAL:O	2.19	0.42
1:A:237:LEU:HD23	1:A:239:TYR:HE2	1.77	0.42
1:A:59:ILE:HA	1:A:64:PHE:O	2.19	0.42
1:F:186:THR:HB	1:F:187:PRO:HD2	2.02	0.42
1:D:177:LYS:HD2	1:D:178:CYS:H	1.84	0.42
1:E:60:TYR:O	1:E:64:PHE:HB2	2.20	0.42
1:E:177:LYS:HB3	1:E:178:CYS:H	1.42	0.42
1:F:236:PHE:O	1:F:239:TYR:N	2.52	0.42
1:D:184:GLY:O	1:D:192:THR:HA	2.20	0.42
1:B:215:LEU:HA	1:B:215:LEU:HD12	1.86	0.42
1:A:146:VAL:HG11	1:A:152:LEU:HD22	2.01	0.42
1:C:237:LEU:HD23	1:C:239:TYR:HE2	1.77	0.42
1:C:146:VAL:HG11	1:C:152:LEU:HD22	2.02	0.42
1:C:136:VAL:HG13	1:C:161:PHE:CD1	2.55	0.42
1:B:177:LYS:HD2	1:B:178:CYS:H	1.85	0.41
1:D:237:LEU:CD2	1:D:239:TYR:HE2	2.34	0.41
1:D:187:PRO:O	1:D:188:ALA:CB	2.68	0.41
1:C:186:THR:HB	1:C:187:PRO:HD2	2.03	0.41
1:B:174:ALA:HB3	1:B:177:LYS:HB2	2.02	0.41
1:E:177:LYS:HE2	2:E:537:HOH:O	2.19	0.41
1:D:234:VAL:O	1:D:239:TYR:HA	2.21	0.41
1:E:146:VAL:HG11	1:E:152:LEU:HD22	2.01	0.41
1:E:59:ILE:HD13	1:E:65:LEU:HD23	2.03	0.41
1:C:140:GLU:OE2	1:C:228:TYR:O	2.38	0.41
1:B:54:ALA:HA	1:B:55:PRO:HD3	1.95	0.41
1:C:221:ARG:C	1:C:252:VAL:HG11	2.41	0.40
1:B:50:LYS:HB2	1:B:53:LYS:HG3	2.03	0.40
1:C:147:GLN:HB3	1:C:147:GLN:HE21	1.67	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:35:ARG:HD2	1:D:35:ARG:HH11	1.71	0.40
1:C:215:LEU:HD12	1:C:215:LEU:HA	1.86	0.40
1:B:35:ARG:HD2	1:B:35:ARG:HH11	1.70	0.40
1:E:136:VAL:HA	1:E:159:SER:O	2.22	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	A	221/276 (80%)	212 (96%)	8 (4%)	1 (0%)	34	48
1	B	221/276 (80%)	212 (96%)	8 (4%)	1 (0%)	34	48
1	C	221/276 (80%)	215 (97%)	6 (3%)	0	100	100
1	D	223/276 (81%)	215 (96%)	8 (4%)	0	100	100
1	E	221/276 (80%)	212 (96%)	9 (4%)	0	100	100
1	F	220/276 (80%)	212 (96%)	8 (4%)	0	100	100
All	All	1327/1656 (80%)	1278 (96%)	47 (4%)	2 (0%)	52	69

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	223	GLU
1	B	95	ALA

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	A	185/212 (87%)	181 (98%)	4 (2%)	60	79
1	B	184/212 (87%)	179 (97%)	5 (3%)	52	73
1	C	184/212 (87%)	180 (98%)	4 (2%)	60	79
1	D	186/212 (88%)	182 (98%)	4 (2%)	60	79
1	E	184/212 (87%)	178 (97%)	6 (3%)	45	66
1	F	183/212 (86%)	180 (98%)	3 (2%)	70	86
All	All	1106/1272 (87%)	1080 (98%)	26 (2%)	57	76

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	83	THR
1	A	106	PRO
1	A	147	GLN
1	A	177	LYS
1	B	35	ARG
1	B	83	THR
1	B	106	PRO
1	B	147	GLN
1	B	177	LYS
1	C	83	THR
1	C	135	GLU
1	C	147	GLN
1	C	177	LYS
1	D	83	THR
1	D	106	PRO
1	D	147	GLN
1	D	177	LYS
1	E	83	THR
1	E	135	GLU
1	E	147	GLN
1	E	151	SER
1	E	177	LYS
1	E	191	TYR
1	F	83	THR
1	F	147	GLN
1	F	177	LYS

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	147	GLN
1	B	147	GLN
1	C	147	GLN
1	D	100	GLN
1	D	147	GLN
1	E	147	GLN
1	F	147	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	225/276 (81%)	-0.39	2 (0%) 85 85	14, 30, 51, 87	0
1	B	225/276 (81%)	-0.37	1 (0%) 93 93	13, 29, 50, 87	0
1	C	225/276 (81%)	-0.23	5 (2%) 65 64	15, 29, 51, 87	0
1	D	227/276 (82%)	-0.25	1 (0%) 93 93	13, 29, 51, 87	0
1	E	225/276 (81%)	-0.19	6 (2%) 58 57	13, 29, 51, 92	0
1	F	224/276 (81%)	-0.21	2 (0%) 85 85	13, 30, 50, 87	0
All	All	1351/1656 (81%)	-0.27	17 (1%) 79 79	13, 29, 51, 92	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	253	SER	3.6
1	E	175	PRO	3.3
1	E	11	GLY	3.2
1	F	175	PRO	3.2
1	E	251	THR	3.0
1	D	253	SER	2.9
1	C	176	GLY	2.7
1	E	222	ALA	2.4
1	A	117	ILE	2.3
1	B	67	SER	2.3
1	C	175	PRO	2.3
1	F	147	GLN	2.3
1	E	176	GLY	2.2
1	C	177	LYS	2.2
1	E	51	PRO	2.1
1	C	11	GLY	2.1
1	C	38	GLN	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.