



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 03:26 am BST

PDB ID : 3I50
Title : Crystal structure of the West Nile Virus envelope glycoprotein in complex with the E53 antibody Fab
Authors : Nybakken, G.E.; Warren, J.T.; Chen, B.R.; Nelson, C.A.; Fremont, D.H.
Deposited on : 2009-07-03
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

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

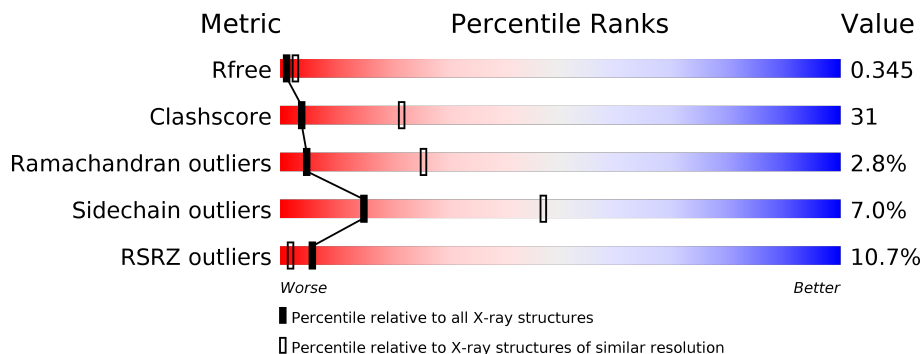
1 Overall quality at a glance

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))
R_{free}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (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 ≥ 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	E	402	 3% 33% 31% 32%
2	L	207	 14% 49% 46%
3	H	221	 14% 45% 48%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5283 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 Envelope glycoprotein.

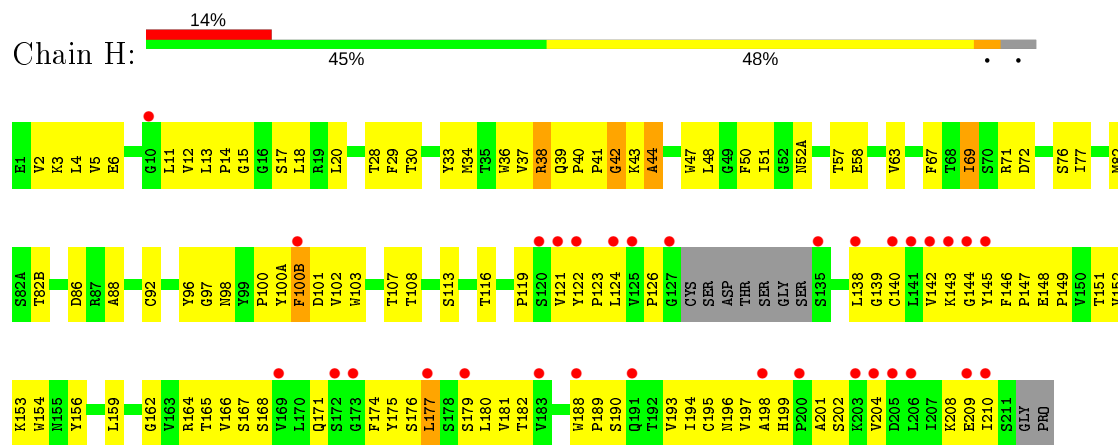
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	E	273	2088	1312	359	399	18	0	0	0

- Molecule 2 is a protein called murine kappa light chain of E53 monoclonal antibody Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	L	207	1583	986	262	326	9	0	0	0

- Molecule 3 is a protein called murine heavy chain (IgG3) of E53 monoclonal antibody Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	H	212	1612	1036	258	312	6	0	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	146.41Å 160.14Å 43.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00 50.15 – 3.00	Depositor EDS
% Data completeness (in resolution range)	93.5 (20.00-3.00) 99.3 (50.15-3.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 3.01Å)	Xtrriage
Refinement program	PHENIX ?	Depositor
R, R_{free}	0.247 , 0.344 0.256 , 0.345	Depositor DCC
R_{free} test set	1097 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	100.7	Xtrriage
Anisotropy	0.893	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 106.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5283	wwPDB-VP
Average B, all atoms (Å ²)	146.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.32% 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 (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	E	0.26	0/2128	0.48	0/2878
2	L	0.29	0/1620	0.49	0/2198
3	H	0.27	0/1654	0.50	0/2262
All	All	0.27	0/5402	0.49	0/7338

There are no bond length outliers.

There are no bond angle outliers.

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	E	2088	0	2057	121	0
2	L	1583	0	1522	104	0
3	H	1612	0	1584	109	0
All	All	5283	0	5163	322	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 31.

All (322) 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:E:79:GLU:HG2	1:E:94:GLN:HE22	1.23	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:168:SER:HB3	3:H:177:LEU:HB3	1.48	0.93
1:E:26:GLU:HG3	1:E:28:ASP:H	1.33	0.90
3:H:51:ILE:HD13	3:H:71:ARG:HG3	1.54	0.89
1:E:42:ASP:HB2	1:E:142:PHE:HB2	1.55	0.88
3:H:148:GLU:HG2	3:H:149:PRO:HA	1.57	0.86
1:E:215:ARG:HH11	1:E:215:ARG:CG	1.90	0.84
1:E:79:GLU:HG2	1:E:94:GLN:NE2	1.93	0.83
1:E:53:LEU:HD22	1:E:130:ILE:HG12	1.59	0.83
2:L:103:LYS:HE3	3:H:43:LYS:HD2	1.58	0.83
1:E:19:THR:HB	1:E:293:GLU:HB2	1.59	0.82
1:E:46:MET:HB2	1:E:138:GLU:HB3	1.63	0.80
1:E:25:LEU:O	1:E:285:HIS:HB2	1.81	0.80
3:H:37:VAL:HG22	3:H:47:TRP:HA	1.62	0.79
1:E:77:MET:HE2	3:H:98:ASN:H	1.48	0.78
1:E:224:PRO:HG2	1:E:239:LEU:HD12	1.65	0.78
2:L:155:ARG:HD3	2:L:179:LEU:HD21	1.65	0.78
1:E:59:TYR:CD2	1:E:225:TRP:HB3	2.18	0.78
3:H:199:HIS:NE2	3:H:201:ALA:HB3	1.99	0.77
1:E:23:LEU:HD11	1:E:31:VAL:HG21	1.66	0.77
3:H:143:LYS:HZ2	3:H:143:LYS:HB3	1.49	0.77
1:E:77:MET:HE2	3:H:98:ASN:N	1.99	0.77
1:E:9:ARG:HG3	1:E:30:CYS:SG	2.24	0.77
2:L:195:GLU:HA	2:L:206:VAL:HG22	1.65	0.77
1:E:15:VAL:HG12	1:E:18:ALA:HB3	1.68	0.75
1:E:248:THR:HG23	1:E:249:LYS:H	1.52	0.74
2:L:122:SER:HA	2:L:125:LEU:HD12	1.70	0.73
2:L:124:GLN:HG2	2:L:129:GLY:HA3	1.70	0.73
2:L:6:ARG:NH2	2:L:88:CYS:H	1.87	0.73
2:L:128:GLY:HA2	2:L:183:LYS:HD2	1.72	0.72
1:E:215:ARG:HH11	1:E:215:ARG:HG3	1.52	0.71
3:H:13:LEU:HD22	3:H:113:SER:HA	1.74	0.70
1:E:44:LYS:HD2	1:E:46:MET:SD	2.32	0.70
1:E:23:LEU:HD21	1:E:31:VAL:HB	1.74	0.69
1:E:65:VAL:HG23	1:E:254:ALA:HB2	1.74	0.69
3:H:208:LYS:HG2	3:H:209:GLU:H	1.57	0.69
2:L:46:ILE:HD13	2:L:47:TRP:N	2.08	0.69
1:E:61:TYR:HB2	1:E:258:GLN:HB2	1.74	0.69
2:L:155:ARG:HH22	2:L:181:LEU:CD1	2.05	0.69
2:L:37:GLN:HG3	2:L:86:TYR:CE1	2.28	0.68
1:E:26:GLU:CG	1:E:29:SER:H	2.06	0.68
2:L:181:LEU:HD23	2:L:182:THR:HG23	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:154:TRP:CZ3	3:H:195:CYS:HB3	2.29	0.68
1:E:26:GLU:HG2	1:E:29:SER:H	1.57	0.68
1:E:30:CYS:HB3	1:E:44:LYS:HG3	1.76	0.67
2:L:155:ARG:NH2	2:L:181:LEU:HB2	2.10	0.67
2:L:46:ILE:HD13	2:L:47:TRP:H	1.59	0.67
1:E:23:LEU:HD12	1:E:24:VAL:H	1.60	0.66
1:E:26:GLU:HG2	1:E:29:SER:HB3	1.77	0.66
1:E:14:GLY:O	1:E:21:VAL:HG22	1.95	0.66
3:H:29:PHE:CD2	3:H:76:SER:HA	2.31	0.66
2:L:7:THR:HB	2:L:22:THR:HB	1.78	0.66
2:L:6:ARG:NH1	2:L:86:TYR:O	2.28	0.66
2:L:125:LEU:HD23	2:L:130:ALA:HB2	1.77	0.65
3:H:3:LYS:HE2	3:H:5:VAL:HG12	1.79	0.65
2:L:78:MET:O	2:L:79:GLU:HG3	1.97	0.65
3:H:123:PRO:HA	3:H:210:ILE:HD12	1.78	0.65
2:L:47:TRP:HZ3	2:L:62:PHE:CE2	2.14	0.64
1:E:201:TYR:HB3	1:E:212:LEU:HD11	1.78	0.64
1:E:58:SER:HB3	1:E:126:ILE:HA	1.78	0.64
2:L:155:ARG:HH22	2:L:181:LEU:HD13	1.61	0.64
2:L:183:LYS:C	2:L:185:GLU:H	2.00	0.64
2:L:183:LYS:C	2:L:183:LYS:HD3	2.19	0.64
2:L:185:GLU:HG3	2:L:188:ARG:HB3	1.79	0.63
3:H:3:LYS:HE2	3:H:5:VAL:CG1	2.29	0.63
3:H:51:ILE:HG23	3:H:71:ARG:HD2	1.80	0.63
3:H:143:LYS:HZ3	3:H:144:GLY:H	1.47	0.62
2:L:47:TRP:CZ3	2:L:62:PHE:CE2	2.86	0.62
1:E:26:GLU:HG2	1:E:29:SER:CB	2.30	0.61
3:H:38:ARG:HG2	3:H:48:LEU:HD21	1.83	0.61
2:L:135:PHE:CE2	3:H:180:LEU:HD11	2.35	0.61
2:L:150:ILE:HB	2:L:153:SER:HB2	1.83	0.61
2:L:185:GLU:HA	2:L:188:ARG:HB3	1.81	0.61
1:E:13:GLU:O	1:E:14:GLY:C	2.39	0.61
3:H:121:VAL:HG22	3:H:142:VAL:HG22	1.83	0.60
2:L:6:ARG:HH21	2:L:99:GLY:HA3	1.67	0.60
1:E:59:TYR:CE2	1:E:225:TRP:HB3	2.36	0.60
3:H:67:PHE:CD1	3:H:82:MET:HA	2.36	0.60
2:L:23:CYS:HB2	2:L:35:TRP:CH2	2.36	0.60
2:L:103:LYS:CE	3:H:43:LYS:HD2	2.31	0.60
3:H:143:LYS:HG2	3:H:176:SER:OG	2.02	0.60
3:H:100(B):PHE:CD1	3:H:100(B):PHE:N	2.69	0.59
2:L:94:HIS:NE2	3:H:58:GLU:HG2	2.16	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:66:GLY:HA3	2:L:71:TYR:CD2	2.37	0.59
2:L:73:LEU:HD12	2:L:73:LEU:C	2.23	0.59
2:L:12:SER:HB3	2:L:107:LYS:HB2	1.85	0.59
1:E:12:LEU:HD13	1:E:31:VAL:HG11	1.85	0.58
2:L:155:ARG:HH21	2:L:181:LEU:HB2	1.68	0.58
2:L:108:ARG:HG3	2:L:109:ALA:O	2.03	0.58
2:L:142:LYS:HB2	2:L:173:TYR:CE2	2.39	0.58
3:H:188:TRP:CD1	3:H:193:VAL:HG21	2.39	0.58
1:E:70:THR:HB	1:E:115:THR:OG1	2.04	0.58
2:L:183:LYS:C	2:L:185:GLU:N	2.58	0.58
1:E:169:ILE:HD12	1:E:174:PRO:O	2.04	0.57
1:E:70:THR:HG21	1:E:250:GLN:H	1.69	0.57
2:L:36:TYR:CE1	2:L:46:ILE:HG12	2.39	0.57
3:H:100(A):TYR:HE1	3:H:101:ASP:HB3	1.69	0.57
3:H:72:ASP:HB3	3:H:77:ILE:HB	1.87	0.57
3:H:100(B):PHE:HD1	3:H:100(B):PHE:N	2.01	0.57
1:E:143:VAL:O	1:E:163:GLN:HG2	2.05	0.57
2:L:6:ARG:HH12	2:L:101:GLY:HA2	1.70	0.57
1:E:223:LEU:O	1:E:225:TRP:HD1	1.87	0.57
1:E:193:ARG:HD2	1:E:287:LYS:NZ	2.20	0.56
2:L:145:ASN:HB2	2:L:197:THR:HG23	1.87	0.56
1:E:14:GLY:O	1:E:15:VAL:HG23	2.05	0.56
3:H:148:GLU:HB2	3:H:175:TYR:CZ	2.40	0.56
2:L:91:TRP:CZ2	3:H:100:PRO:HG2	2.41	0.56
3:H:40:PRO:HD2	3:H:44:ALA:O	2.06	0.56
2:L:37:GLN:HG3	2:L:86:TYR:CZ	2.41	0.56
2:L:47:TRP:HZ3	2:L:62:PHE:CZ	2.22	0.56
1:E:272:VAL:HG12	1:E:273:GLU:N	2.20	0.56
1:E:53:LEU:HD11	1:E:279:VAL:HG21	1.88	0.56
3:H:179:SER:C	3:H:180:LEU:HD12	2.26	0.56
3:H:33:TYR:O	3:H:34:MET:HG2	2.06	0.56
2:L:47:TRP:CZ3	2:L:62:PHE:CZ	2.94	0.56
2:L:175:MET:HG2	2:L:176:SER:N	2.21	0.55
3:H:152:VAL:HA	3:H:196:ASN:O	2.06	0.55
3:H:199:HIS:CD2	3:H:202:SER:H	2.25	0.55
2:L:107:LYS:HA	2:L:140:TYR:OH	2.07	0.55
2:L:198:HIS:CG	2:L:199:LYS:H	2.25	0.55
1:E:96:VAL:HG13	1:E:110:LYS:HD2	1.88	0.54
3:H:152:VAL:HG22	3:H:197:VAL:HG13	1.88	0.54
2:L:150:ILE:HD11	2:L:155:ARG:HD2	1.88	0.54
1:E:215:ARG:HH11	1:E:215:ARG:HG2	1.70	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:6:ARG:HH22	2:L:87:TYR:HA	1.72	0.54
1:E:200:ALA:O	1:E:215:ARG:HB2	2.08	0.54
3:H:188:TRP:CD1	3:H:189:PRO:HA	2.42	0.54
2:L:179:LEU:HD12	2:L:180:THR:H	1.73	0.54
3:H:41:PRO:HD3	3:H:88:ALA:HA	1.89	0.54
2:L:35:TRP:HB2	2:L:48:ILE:HB	1.90	0.54
1:E:15:VAL:CG1	1:E:18:ALA:HB3	2.37	0.53
3:H:63:VAL:HG12	3:H:63:VAL:O	2.06	0.53
1:E:64:THR:OG1	1:E:120:ALA:HB3	2.09	0.53
2:L:128:GLY:C	2:L:183:LYS:HB3	2.28	0.53
3:H:101:ASP:O	3:H:103:TRP:HD1	1.90	0.53
2:L:151:ASP:OD2	2:L:189:HIS:CD2	2.60	0.53
1:E:175:SER:O	1:E:176:TYR:HB2	2.08	0.53
1:E:53:LEU:CD2	1:E:130:ILE:HG12	2.37	0.53
1:E:19:THR:HB	1:E:293:GLU:CB	2.34	0.53
2:L:73:LEU:HD12	2:L:73:LEU:O	2.07	0.53
3:H:100(A):TYR:CE1	3:H:101:ASP:HB3	2.44	0.53
3:H:15:GLY:O	3:H:82(B):THR:HA	2.09	0.52
3:H:156:TYR:H	3:H:196:ASN:HD21	1.57	0.52
3:H:29:PHE:CG	3:H:76:SER:HA	2.45	0.52
1:E:269:ALA:O	1:E:271:PRO:HD3	2.09	0.52
1:E:7:SER:C	1:E:9:ARG:H	2.13	0.52
2:L:140:TYR:CG	2:L:141:PRO:HA	2.45	0.52
3:H:177:LEU:HD12	3:H:177:LEU:O	2.09	0.52
2:L:47:TRP:HZ3	2:L:62:PHE:CD2	2.28	0.52
1:E:187:THR:HB	1:E:291:LYS:HB2	1.90	0.52
1:E:172:ALA:O	1:E:174:PRO:HD3	2.10	0.51
1:E:291:LYS:HB3	1:E:294:LYS:HE2	1.92	0.51
2:L:91:TRP:CE2	3:H:100:PRO:HD2	2.45	0.51
3:H:36:TRP:CD1	3:H:69:ILE:HD12	2.46	0.51
2:L:123:GLU:HB3	3:H:122:TYR:CE1	2.45	0.51
1:E:93:ARG:HG3	1:E:242:PHE:CE1	2.46	0.51
1:E:270:ILE:O	1:E:272:VAL:HG23	2.11	0.51
3:H:36:TRP:HD1	3:H:69:ILE:HD12	1.76	0.51
3:H:116:THR:HG23	3:H:201:ALA:HB1	1.92	0.50
1:E:193:ARG:HD2	1:E:287:LYS:HZ3	1.75	0.50
2:L:117:ILE:HD12	2:L:194:CYS:SG	2.51	0.50
1:E:224:PRO:HA	1:E:234:ARG:O	2.11	0.50
2:L:150:ILE:CD1	2:L:155:ARG:HD2	2.40	0.50
3:H:124:LEU:HB2	3:H:139:GLY:O	2.11	0.50
1:E:131:LEU:HD21	1:E:199:ASN:HB2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:103:TRP:N	3:H:103:TRP:CD1	2.79	0.50
3:H:202:SER:O	3:H:204:VAL:HG13	2.12	0.50
1:E:65:VAL:HG22	1:E:240:MET:HE2	1.94	0.50
2:L:137:ASN:HD22	3:H:164:ARG:HG3	1.77	0.49
2:L:150:ILE:HG12	2:L:155:ARG:HD2	1.94	0.49
3:H:153:LYS:HB2	3:H:196:ASN:HB2	1.94	0.49
3:H:50:PHE:O	3:H:57:THR:HG23	2.12	0.49
3:H:143:LYS:NZ	3:H:144:GLY:H	2.10	0.49
3:H:188:TRP:HA	3:H:190:SER:H	1.77	0.49
1:E:96:VAL:CG1	1:E:110:LYS:HD2	2.42	0.49
3:H:28:THR:O	3:H:28:THR:HG22	2.11	0.49
1:E:200:ALA:O	1:E:201:TYR:HD1	1.95	0.49
1:E:24:VAL:O	1:E:25:LEU:HD23	2.12	0.49
3:H:194:ILE:HG22	3:H:209:GLU:HB3	1.95	0.49
2:L:155:ARG:NH2	2:L:181:LEU:CB	2.76	0.49
1:E:216:GLU:O	1:E:220:ASP:HB2	2.13	0.48
1:E:15:VAL:HB	1:E:20:TRP:O	2.13	0.48
1:E:180:LEU:HB2	1:E:184:GLY:O	2.13	0.48
1:E:202:TYR:O	1:E:212:LEU:HD12	2.13	0.48
1:E:246:HIS:O	1:E:247:ALA:C	2.52	0.48
1:E:26:GLU:HG3	1:E:27:GLY:N	2.29	0.48
3:H:147:PRO:HD2	3:H:201:ALA:HB1	1.96	0.47
1:E:196:ILE:HG12	1:E:197:ASP:N	2.29	0.47
2:L:124:GLN:C	2:L:126:THR:H	2.16	0.47
2:L:2:ILE:O	2:L:97:THR:HG21	2.14	0.47
3:H:17:SER:HB2	3:H:82:MET:O	2.14	0.47
3:H:38:ARG:NH2	3:H:86:ASP:HA	2.30	0.47
1:E:23:LEU:HD21	1:E:31:VAL:CB	2.44	0.47
2:L:108:ARG:HG3	2:L:109:ALA:N	2.29	0.47
1:E:214:HIS:O	1:E:215:ARG:C	2.54	0.47
1:E:40:THR:OG1	1:E:144:HIS:ND1	2.47	0.47
2:L:73:LEU:CD1	2:L:73:LEU:C	2.83	0.47
1:E:57:ARG:HD2	1:E:225:TRP:CE3	2.50	0.47
3:H:143:LYS:NZ	3:H:143:LYS:HB3	2.24	0.47
3:H:171:GLN:O	3:H:174:PHE:HB2	2.14	0.46
2:L:14:SER:O	2:L:17:GLU:HB3	2.15	0.46
2:L:15:PRO:HG3	2:L:106:LEU:HD11	1.97	0.46
3:H:139:GLY:HA3	3:H:179:SER:O	2.15	0.46
3:H:96:TYR:O	3:H:100:PRO:HA	2.16	0.46
1:E:24:VAL:HA	1:E:286:LEU:O	2.15	0.46
1:E:84:LYS:HD3	1:E:84:LYS:HA	1.71	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:148:GLU:CG	3:H:149:PRO:HA	2.36	0.46
3:H:2:VAL:CG1	3:H:102:VAL:HG11	2.46	0.46
3:H:101:ASP:O	3:H:103:TRP:CD1	2.68	0.46
1:E:215:ARG:NH1	1:E:215:ARG:CG	2.61	0.46
1:E:258:GLN:O	1:E:259:GLU:C	2.53	0.46
3:H:67:PHE:HD1	3:H:82:MET:HA	1.78	0.46
1:E:10:ASP:HB2	1:E:31:VAL:HG22	1.96	0.46
1:E:57:ARG:O	1:E:218:PHE:HZ	1.99	0.46
2:L:150:ILE:CG1	2:L:155:ARG:HD2	2.46	0.46
2:L:26:SER:O	2:L:27:SER:HB2	2.16	0.46
2:L:7:THR:HB	2:L:22:THR:CB	2.44	0.46
3:H:12:VAL:HG21	3:H:18:LEU:HD22	1.98	0.45
3:H:199:HIS:CE1	3:H:201:ALA:HB3	2.49	0.45
1:E:119:PHE:CG	1:E:236:ARG:HD2	2.50	0.45
3:H:51:ILE:HG23	3:H:51:ILE:O	2.15	0.45
1:E:183:TYR:O	1:E:296:GLN:HB3	2.16	0.45
3:H:38:ARG:HH21	3:H:86:ASP:HA	1.81	0.45
2:L:130:ALA:N	2:L:183:LYS:HB2	2.31	0.45
1:E:50:ALA:HB1	1:E:53:LEU:HD21	1.99	0.45
1:E:84:LYS:C	1:E:86:ALA:H	2.20	0.45
1:E:80:ALA:H	1:E:112:SER:CB	2.30	0.45
3:H:166:VAL:HG23	3:H:167:SER:O	2.15	0.45
2:L:36:TYR:CD1	2:L:46:ILE:HA	2.51	0.45
1:E:193:ARG:NH1	1:E:287:LYS:HE2	2.32	0.45
1:E:295:LEU:HD12	1:E:296:GLN:H	1.82	0.45
1:E:53:LEU:HD11	1:E:279:VAL:CG2	2.46	0.45
3:H:6:GLU:HB3	3:H:107:THR:HB	1.99	0.45
1:E:285:HIS:NE2	1:E:287:LYS:HD2	2.33	0.44
1:E:41:ILE:HG22	1:E:143:VAL:HG22	1.99	0.44
2:L:166:GLN:HG3	2:L:171:SER:O	2.17	0.44
3:H:154:TRP:CH2	3:H:195:CYS:HB3	2.52	0.44
1:E:183:TYR:O	1:E:295:LEU:HD12	2.17	0.44
1:E:48:MET:HE2	1:E:281:LEU:HB2	2.00	0.44
3:H:199:HIS:HD2	3:H:202:SER:OG	2.01	0.44
3:H:63:VAL:HG13	3:H:67:PHE:CD2	2.53	0.44
3:H:181:VAL:HG22	3:H:182:THR:N	2.33	0.44
2:L:38:GLN:O	2:L:84:ALA:HB1	2.18	0.44
1:E:137:TYR:CE2	1:E:192:PRO:HB3	2.52	0.44
2:L:194:CYS:O	2:L:206:VAL:HA	2.18	0.44
1:E:202:TYR:HE1	1:E:215:ARG:HD2	1.83	0.43
3:H:121:VAL:HG12	3:H:208:LYS:HD2	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:14:SER:O	2:L:17:GLU:CB	2.66	0.43
1:E:50:ALA:CB	1:E:130:ILE:HD13	2.47	0.43
1:E:215:ARG:HG3	1:E:215:ARG:NH1	2.26	0.43
1:E:50:ALA:HB1	1:E:130:ILE:HG21	2.00	0.43
1:E:74:CYS:HB3	1:E:75:PRO:HD2	2.00	0.43
2:L:42:THR:HG22	2:L:43:SER:O	2.18	0.43
3:H:165:THR:HA	3:H:179:SER:HA	2.00	0.43
3:H:30:THR:O	3:H:52(A):ASN:HB2	2.18	0.43
1:E:200:ALA:C	1:E:201:TYR:HD1	2.21	0.43
2:L:137:ASN:O	2:L:138:ASN:C	2.56	0.43
1:E:12:LEU:HD13	1:E:31:VAL:CG1	2.48	0.43
1:E:93:ARG:HG3	1:E:242:PHE:CZ	2.54	0.43
3:H:140:CYS:SG	3:H:210:ILE:HD11	2.59	0.43
2:L:10:ILE:HG12	2:L:11:MET:N	2.34	0.43
1:E:227:SER:H	1:E:230:SER:HB3	1.82	0.43
3:H:63:VAL:CG1	3:H:63:VAL:O	2.67	0.42
3:H:71:ARG:HB3	3:H:71:ARG:HE	1.68	0.42
2:L:151:ASP:N	2:L:151:ASP:OD1	2.52	0.42
2:L:85:THR:HG23	2:L:103:LYS:HD3	2.00	0.42
1:E:195:GLY:HA3	1:E:283:SER:HB2	2.00	0.42
1:E:47:ASN:O	1:E:138:GLU:N	2.51	0.42
3:H:119:PRO:HG3	3:H:145:TYR:CB	2.49	0.42
3:H:96:TYR:CD2	3:H:97:GLY:N	2.87	0.42
1:E:246:HIS:O	1:E:248:THR:N	2.53	0.42
3:H:2:VAL:HG11	3:H:102:VAL:HG11	2.01	0.42
2:L:6:ARG:NH2	2:L:87:TYR:HA	2.33	0.42
2:L:197:THR:HB	2:L:204:PRO:HB3	2.00	0.42
1:E:13:GLU:O	1:E:13:GLU:HG3	2.19	0.42
3:H:168:SER:HA	3:H:176:SER:O	2.19	0.42
1:E:175:SER:HA	1:E:188:VAL:O	2.19	0.42
3:H:188:TRP:CA	3:H:190:SER:H	2.32	0.42
2:L:138:ASN:HD21	3:H:164:ARG:HH11	1.68	0.42
2:L:129:GLY:C	2:L:183:LYS:HG3	2.40	0.42
1:E:197:ASP:O	1:E:200:ALA:HB3	2.18	0.42
2:L:80:ALA:O	2:L:83:VAL:HG23	2.19	0.42
1:E:193:ARG:HH11	1:E:287:LYS:HE2	1.85	0.41
2:L:135:PHE:CD2	3:H:180:LEU:HD11	2.54	0.41
2:L:135:PHE:C	2:L:136:LEU:HD12	2.41	0.41
1:E:68:LEU:HD12	1:E:68:LEU:HA	1.83	0.41
1:E:70:THR:HG21	1:E:250:GLN:N	2.33	0.41
1:E:99:ARG:NH1	1:E:103:ASN:O	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:47:TRP:CZ3	2:L:62:PHE:CD2	3.08	0.41
2:L:96:LEU:HD23	2:L:96:LEU:N	2.34	0.41
2:L:134:CYS:HB2	2:L:148:TRP:CH2	2.55	0.41
2:L:30:VAL:CG1	2:L:31:SER:N	2.84	0.41
2:L:115:VAL:HA	2:L:135:PHE:O	2.20	0.41
1:E:171:PRO:HA	1:E:192:PRO:HG2	2.02	0.41
1:E:83:ASP:OD1	1:E:83:ASP:N	2.54	0.41
3:H:116:THR:HA	3:H:146:PHE:O	2.21	0.41
2:L:142:LYS:O	2:L:144:ILE:HG22	2.21	0.41
2:L:151:ASP:CG	2:L:191:SER:HB2	2.40	0.41
1:E:169:ILE:HG23	1:E:174:PRO:HA	2.02	0.41
2:L:146:VAL:HG21	2:L:175:MET:SD	2.61	0.41
1:E:87:ASP:HA	1:E:88:PRO:HD3	1.94	0.41
3:H:11:LEU:HB2	3:H:147:PRO:HG3	2.03	0.41
2:L:50:GLU:O	2:L:51:SER:HB2	2.21	0.41
3:H:142:VAL:HB	3:H:177:LEU:HD12	2.03	0.41
3:H:145:TYR:CZ	3:H:175:TYR:HB2	2.56	0.41
3:H:3:LYS:HE2	3:H:5:VAL:HG11	2.03	0.41
3:H:42:GLY:O	3:H:43:LYS:HB2	2.21	0.41
2:L:6:ARG:NH1	2:L:101:GLY:HA2	2.35	0.41
3:H:126:PRO:HD3	3:H:138:LEU:HG	2.03	0.41
2:L:34:HIS:HD2	2:L:49:TYR:HA	1.86	0.41
3:H:138:LEU:HB3	3:H:139:GLY:H	1.72	0.40
3:H:164:ARG:N	3:H:180:LEU:O	2.42	0.40
3:H:151:THR:OG1	3:H:198:ALA:HB3	2.21	0.40
2:L:34:HIS:CD2	2:L:49:TYR:HA	2.56	0.40
3:H:162:GLY:O	3:H:181:VAL:HA	2.21	0.40
2:L:6:ARG:NE	2:L:88:CYS:SG	2.95	0.40
1:E:69:SER:O	1:E:82:ASN:ND2	2.54	0.40
3:H:188:TRP:HA	3:H:190:SER:N	2.36	0.40
3:H:39:GLN:O	3:H:88:ALA:HB1	2.21	0.40
1:E:236:ARG:O	1:E:237:GLU:C	2.59	0.40
1:E:221:LEU:HD23	1:E:221:LEU:HA	1.97	0.40
2:L:6:ARG:NH2	2:L:99:GLY:HA3	2.34	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	E	269/402 (67%)	207 (77%)	51 (19%)	11 (4%)	3	16
2	L	205/207 (99%)	164 (80%)	36 (18%)	5 (2%)	6	29
3	H	208/221 (94%)	174 (84%)	31 (15%)	3 (1%)	11	43
All	All	682/830 (82%)	545 (80%)	118 (17%)	19 (3%)	5	25

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	16	SER
1	E	215	ARG
1	E	14	GLY
1	E	36	LYS
1	E	52	ASN
1	E	69	SER
1	E	247	ALA
2	L	27	SER
2	L	44	PRO
2	L	138	ASN
1	E	176	TYR
1	E	248	THR
2	L	174	SER
3	H	42	GLY
3	H	44	ALA
1	E	8	ASN
3	H	14	PRO
2	L	113	PRO
1	E	74	CYS

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	E	226/332 (68%)	207 (92%)	19 (8%)	11	38
2	L	182/182 (100%)	169 (93%)	13 (7%)	14	46
3	H	180/187 (96%)	171 (95%)	9 (5%)	24	60
All	All	588/701 (84%)	547 (93%)	41 (7%)	15	47

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	15	VAL
1	E	22	ASP
1	E	30	CYS
1	E	37	ASP
1	E	41	ILE
1	E	58	SER
1	E	67	ASP
1	E	70	THR
1	E	103	ASN
1	E	113	ILE
1	E	126	ILE
1	E	130	ILE
1	E	178	LEU
1	E	186	VAL
1	E	196	ILE
1	E	198	THR
1	E	215	ARG
1	E	249	LYS
1	E	281	LEU
2	L	22	THR
2	L	46	ILE
2	L	69	THR
2	L	73	LEU
2	L	76	SER
2	L	92	SER
2	L	97	THR

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Mol	Chain	Res	Type
2	L	123	GLU
2	L	134	CYS
2	L	138	ASN
2	L	151	ASP
2	L	189	HIS
2	L	193	THR
3	H	4	LEU
3	H	20	LEU
3	H	38	ARG
3	H	69	ILE
3	H	92	CYS
3	H	100(B)	PHE
3	H	108	THR
3	H	159	LEU
3	H	177	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	94	GLN
2	L	1	GLN
2	L	34	HIS
2	L	137	ASN
3	H	196	ASN
3	H	199	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](#)

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	E	273/402 (67%)	0.38	12 (4%) 34 13	70, 130, 191, 236	0
2	L	207/207 (100%)	0.88	30 (14%) 2 1	66, 134, 229, 275	0
3	H	212/221 (95%)	0.88	32 (15%) 2 1	67, 148, 276, 316	0
All	All	692/830 (83%)	0.68	74 (10%) 6 2	66, 135, 244, 316	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	120	PRO	12.4
2	L	155	ARG	10.3
3	H	127	GLY	9.8
3	H	138	LEU	7.7
3	H	198	ALA	6.9
3	H	205	ASP	6.2
2	L	194	CYS	5.6
3	H	121	VAL	5.0
1	E	11	PHE	5.0
2	L	121	SER	5.0
3	H	142	VAL	4.9
3	H	183	VAL	4.9
2	L	208	SER	4.5
2	L	192	TYR	4.1
1	E	297	LEU	3.9
2	L	127	SER	3.9
2	L	195	GLU	3.9
1	E	165	GLY	3.8
2	L	132	VAL	3.8
2	L	156	GLN	3.8
3	H	120	SER	3.7
3	H	145	TYR	3.7
3	H	177	LEU	3.6

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Mol	Chain	Res	Type	RSRZ
2	L	115	VAL	3.6
2	L	126	THR	3.6
3	H	200	PRO	3.6
1	E	20	TRP	3.4
3	H	143	LYS	3.3
2	L	125	LEU	3.2
3	H	122	TYR	3.2
3	H	172	SER	3.2
2	L	159	VAL	3.2
2	L	122	SER	3.1
3	H	135	SER	3.1
3	H	125	VAL	3.1
3	H	206	LEU	3.0
2	L	196	ALA	3.0
2	L	197	THR	2.9
2	L	133	VAL	2.8
3	H	144	GLY	2.8
2	L	147	LYS	2.7
3	H	10	GLY	2.7
3	H	173	GLY	2.7
1	E	290	VAL	2.6
3	H	141	LEU	2.6
2	L	136	LEU	2.6
3	H	204	VAL	2.6
2	L	91	TRP	2.5
1	E	16	SER	2.5
3	H	169	VAL	2.5
2	L	179	LEU	2.4
3	H	209	GLU	2.3
3	H	210	ILE	2.3
3	H	124	LEU	2.3
1	E	206	VAL	2.3
3	H	188	TRP	2.3
1	E	204	MET	2.3
1	E	196	ILE	2.3
3	H	179	SER	2.3
2	L	4	LEU	2.2
3	H	203	LYS	2.2
3	H	140	CYS	2.2
1	E	7	SER	2.2
2	L	206	VAL	2.2
3	H	100(B)	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
2	L	2	ILE	2.2
2	L	184	ASP	2.1
2	L	119	PRO	2.1
1	E	41	ILE	2.1
2	L	116	SER	2.1
1	E	141	ILE	2.0
3	H	191	GLN	2.0
2	L	134	CYS	2.0
2	L	160	LEU	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.