



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:10 AM UTC

PDB ID : 3I50 / pdb_00003i50
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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

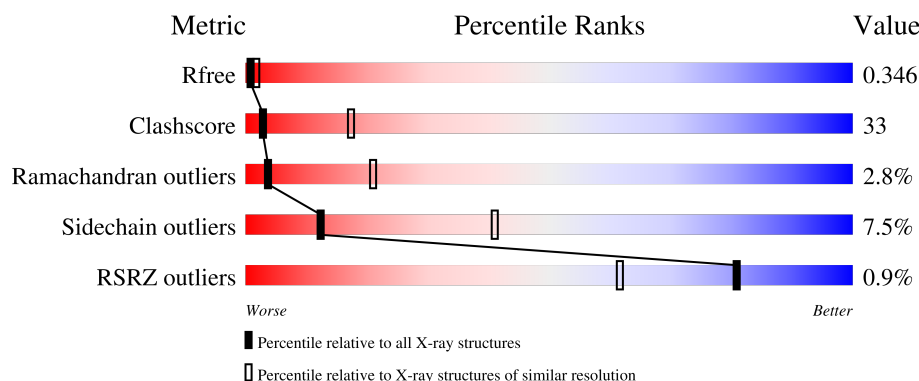
1 Overall quality at a glance

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}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	
2	L	207	
3	H	221	

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
1	E	273	Total	C	N	O	S	0	0	0
			2088	1312	359	399	18			

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

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

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

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



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 20.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	93.5 (20.00-3.00) 99.0 (20.00-3.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 3.01Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.247 , 0.344 0.261 , 0.346	Depositor DCC
R_{free} test set	1098 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	100.7	Xtriage
Anisotropy	0.893	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 106.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5283	wwPDB-VP
Average B, all atoms (Å ²)	146.0	wwPDB-VP

Xtriage'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	0.33	0/2128	0.78	0/2878
2	L	0.36	0/1620	0.78	0/2198
3	H	0.33	0/1654	0.81	2/2262 (0.1%)
All	All	0.34	0/5402	0.79	2/7338 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	118	ALA	CA-C-N	5.60	126.38	120.11
3	H	118	ALA	C-N-CA	5.60	126.38	120.11

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	129	0
2	L	1583	0	1522	108	0
3	H	1612	0	1584	115	0
All	All	5283	0	5163	340	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 33.

All (340) 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
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	1.47	0.79
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:HB3	3:H:143:LYS:HZ2	1.49	0.77
1:E:9:ARG:HG3	1:E:30:CYS:SG	2.24	0.77
1:E:77:MET:HE2	3:H:98:ASN:N	1.99	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:48:MET:HE3	1:E:281:LEU:HB2	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
1:E:61:TYR:HB2	1:E:258:GLN:HB2	1.74	0.69
2:L:46:ILE:HD13	2:L:47:TRP:N	2.08	0.69
2:L:155:ARG:HH22	2:L:181:LEU:CD1	2.05	0.69
2:L:183:LYS:C	2:L:185:GLU:H	2.00	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:37:GLN:HG3	2:L:86:TYR:CE1	2.28	0.68
2:L:183:LYS:C	2:L:183:LYS:HD3	2.19	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
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:6:ARG:NH1	2:L:86:TYR:O	2.28	0.66
2:L:7:THR:HB	2:L:22:THR:HB	1.78	0.66
1:E:13:GLU:O	1:E:14:GLY:C	2.39	0.65
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:73:LEU:C	2:L:73:LEU:HD12	2.23	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
2:L:47:TRP:CZ3	2:L:62:PHE:CE2	2.86	0.62
3:H:143:LYS:HZ3	3:H:144:GLY:H	1.47	0.62
1:E:26:GLU:HG2	1:E:29:SER:CB	2.30	0.61
2:L:135:PHE:CE2	3:H:180:LEU:HD11	2.35	0.61
3:H:38:ARG:HG2	3:H:48:LEU:HD21	1.83	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
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
2:L:23:CYS:HB2	2:L:35:TRP:CH2	2.36	0.60
3:H:67:PHE:CD1	3:H:82:MET:HA	2.36	0.60
3:H:179:SER:C	3:H:180:LEU:HD12	2.26	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
2:L:183:LYS:C	2:L:185:GLU:N	2.58	0.59
2:L:94:HIS:NE2	3:H:58:GLU:HG2	2.16	0.59
3:H:100(B):PHE:CD1	3:H:100(B):PHE:N	2.69	0.59
2:L:66:GLY:HA3	2:L:71:TYR:CD2	2.37	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:128:GLY:C	2:L:183:LYS:HB3	2.28	0.58
1:E:70:THR:HG21	1:E:250:GLN:H	1.69	0.57
1:E:169:ILE:HD12	1:E:174:PRO:O	2.04	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:N	3:H:100(B):PHE:HD1	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
1:E:7:SER:C	1:E:9:ARG:H	2.13	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:37:GLN:HG3	2:L:86:TYR:CZ	2.41	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:47:TRP:HZ3	2:L:62:PHE:CZ	2.22	0.56
1:E:53:LEU:HD11	1:E:279:VAL:HG21	1.88	0.56
1:E:272:VAL:HG12	1:E:273:GLU:N	2.20	0.56
2:L:47:TRP:CZ3	2:L:62:PHE:CZ	2.94	0.56
3:H:33:TYR:O	3:H:34:MET:HG2	2.06	0.56
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:150:ILE:HD11	2:L:155:ARG:HD2	1.88	0.54
3:H:152:VAL:HG22	3:H:197:VAL:HG13	1.88	0.54
1:E:215:ARG:HH11	1:E:215:ARG:HG2	1.70	0.54
2:L:6:ARG:HH22	2:L:87:TYR:HA	1.72	0.54
2:L:163:TRP:CE2	2:L:175:MET:HE3	2.42	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
2:L:35:TRP:HB2	2:L:48:ILE:HB	1.90	0.54
3:H:41:PRO:HD3	3:H:88:ALA:HA	1.89	0.54
2:L:175:MET:HG2	2:L:176:SER:N	2.21	0.53
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:151:ASP:OD2	2:L:189:HIS:CD2	2.60	0.53
3:H:101:ASP:O	3:H:103:TRP:HD1	1.90	0.53
1:E:175:SER:O	1:E:176:TYR:HB2	2.08	0.53
1:E:19:THR:HB	1:E:293:GLU:CB	2.34	0.53
1:E:53:LEU:CD2	1:E:130:ILE:HG12	2.37	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
1:E:246:HIS:O	1:E:247:ALA:C	2.52	0.53
3:H:15:GLY:O	3:H:82(B):THR:HA	2.09	0.52
3:H:29:PHE:CG	3:H:76:SER:HA	2.45	0.52
3:H:156:TYR:H	3:H:196:ASN:HD21	1.57	0.52
1:E:269:ALA:O	1:E:271:PRO:HD3	2.09	0.52
2:L:140:TYR:CG	2:L:141:PRO:HA	2.45	0.52
2:L:47:TRP:HZ3	2:L:62:PHE:CD2	2.28	0.52
3:H:177:LEU:O	3:H:177:LEU:HD12	2.09	0.52
1:E:187:THR:HB	1:E:291:LYS:HB2	1.90	0.52
2:L:73:LEU:C	2:L:73:LEU:CD1	2.83	0.52
2:L:124:GLN:C	2:L:126:THR:H	2.16	0.52
1:E:172:ALA:O	1:E:174:PRO:HD3	2.10	0.51
1:E:214:HIS:O	1:E:215:ARG:C	2.54	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:258:GLN:O	1:E:259:GLU:C	2.53	0.51
1:E:270:ILE:O	1:E:272:VAL:HG23	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
3:H:103:TRP:CD1	3:H:103:TRP:N	2.79	0.50
3:H:202:SER:O	3:H:204:VAL:HG13	2.12	0.50
1:E:204:MET:HE2	1:E:213:VAL:HG11	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
1:E:84:LYS:C	1:E:86:ALA:H	2.20	0.49
3:H:50:PHE:O	3:H:57:THR:HG23	2.12	0.49
3:H:153:LYS:HB2	3:H:196:ASN:HB2	1.94	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
2:L:155:ARG:NH2	2:L:181:LEU:CB	2.76	0.49
3:H:194:ILE:HG22	3:H:209:GLU:HB3	1.95	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:200:ALA:C	1:E:201:TYR:HD1	2.21	0.48
1:E:26:GLU:HG3	1:E:27:GLY:N	2.29	0.48
1:E:65:VAL:HG22	1:E:240:MET:HE2	1.96	0.48
2:L:137:ASN:O	2:L:138:ASN:C	2.56	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:2:ILE:O	2:L:97:THR:HG21	2.14	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
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:40:THR:OG1	1:E:144:HIS:ND1	2.47	0.47
1:E:57:ARG:HD2	1:E:225:TRP:CE3	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:143:LYS:HB3	3:H:143:LYS:NZ	2.24	0.47
2:L:151:ASP:CG	2:L:191:SER:HB2	2.40	0.46
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
2:L:129:GLY:C	2:L:183:LYS:HG3	2.40	0.46
2:L:135:PHE:C	2:L:136:LEU:HD12	2.41	0.46
3:H:96:TYR:O	3:H:100:PRO:HA	2.16	0.46
3:H:139:GLY:HA3	3:H:179:SER:O	2.15	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
3:H:2:VAL:CG1	3:H:102:VAL:HG11	2.46	0.46
3:H:148:GLU:CG	3:H:149:PRO:HA	2.36	0.46
3:H:101:ASP:O	3:H:103:TRP:CD1	2.68	0.46
1:E:215:ARG:CG	1:E:215:ARG:NH1	2.61	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:7:THR:HB	2:L:22:THR:CB	2.44	0.46
2:L:26:SER:O	2:L:27:SER:HB2	2.16	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:48:MET:HE1	1:E:283:SER:C	2.41	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
2:L:130:ALA:N	2:L:183:LYS:HB2	2.31	0.45
3:H:38:ARG:HH21	3:H:86:ASP:HA	1.81	0.45
1:E:50:ALA:HB1	1:E:53:LEU:HD21	1.99	0.45
1:E:236:ARG:O	1:E:237:GLU:C	2.59	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:53:LEU:HD11	1:E:279:VAL:CG2	2.46	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
3:H:6:GLU:HB3	3:H:107:THR:HB	1.99	0.45
1:E:41:ILE:HG22	1:E:143:VAL:HG22	1.99	0.44
1:E:285:HIS:NE2	1:E:287:LYS:HD2	2.33	0.44
2:L:166:GLN:HG3	2:L:171:SER:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:43:LYS:O	3:H:44:ALA:C	2.59	0.44
2:L:161:ASN:HB3	2:L:175:MET:HE2	1.99	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
2:L:39:LYS:O	2:L:40:SER:C	2.61	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
2:L:38:GLN:O	2:L:84:ALA:HB1	2.18	0.44
3:H:181:VAL:HG22	3:H:182:THR:N	2.33	0.44
1:E:137:TYR:CE2	1:E:192:PRO:HB3	2.52	0.44
1:E:260:GLY:C	1:E:262:LEU:H	2.26	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
2:L:14:SER:O	2:L:17:GLU:CB	2.66	0.43
3:H:121:VAL:HG12	3:H:208:LYS:HD2	1.99	0.43
1:E:50:ALA:HB1	1:E:130:ILE:HG21	2.00	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: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:30:THR:O	3:H:52(A):ASN:HB2	2.18	0.43
3:H:165:THR:HA	3:H:179:SER:HA	2.00	0.43
2:L:124:GLN:C	2:L:126:THR:N	2.75	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
1:E:74:CYS:C	1:E:75:PRO:O	2.61	0.43
2:L:10:ILE:HG12	2:L:11:MET:N	2.34	0.43
3:H:140:CYS:SG	3:H:210:ILE:HD11	2.59	0.43
1:E:227:SER:H	1:E:230:SER:HB3	1.82	0.43
2:L:85:THR:HG23	2:L:103:LYS:HD3	2.00	0.42
2:L:151:ASP:OD1	2:L:151:ASP:N	2.52	0.42
3:H:63:VAL:O	3:H:63:VAL:CG1	2.67	0.42
3:H:71:ARG:HE	3:H:71:ARG:HB3	1.68	0.42
1:E:47:ASN:O	1:E:138:GLU:N	2.51	0.42
1:E:51:ALA:O	1:E:52:ASN:C	2.61	0.42
1:E:195:GLY:HA3	1:E:283:SER:HB2	2.00	0.42
3:H:96:TYR:CD2	3:H:97:GLY:N	2.87	0.42
3:H:119:PRO:HG3	3:H:145:TYR:CB	2.49	0.42
1:E:246:HIS:O	1:E:248:THR:N	2.53	0.42
2:L:6:ARG:NH2	2:L:87:TYR:HA	2.33	0.42
3:H:2:VAL:HG11	3:H:102:VAL:HG11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
3:H:177:LEU:HD12	3:H:177:LEU:C	2.44	0.42
1:E:175:SER:HA	1:E:188:VAL:O	2.19	0.42
2:L:138:ASN:HD21	3:H:164:ARG:HH11	1.68	0.42
3:H:188:TRP:CA	3:H:190:SER:H	2.32	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
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
2:L:47:TRP:CZ3	2:L:62:PHE:CD2	3.08	0.41
2:L:96:LEU:N	2:L:96:LEU:HD23	2.34	0.41
2:L:134:CYS:HB2	2:L:148:TRP:CH2	2.55	0.41
3:H:184:PRO:C	3:H:186:SER:H	2.29	0.41
2:L:115:VAL:HA	2:L:135:PHE:O	2.20	0.41
1:E:83:ASP:OD1	1:E:83:ASP:N	2.54	0.41
1:E:84:LYS:C	1:E:86:ALA:N	2.77	0.41
1:E:171:PRO:HA	1:E:192:PRO:HG2	2.02	0.41
2:L:142:LYS:O	2:L:144:ILE:HG22	2.21	0.41
3:H:116:THR:HA	3:H:146:PHE:O	2.21	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
2:L:6:ARG:NH1	2:L:101:GLY:HA2	2.35	0.41
2:L:122:SER:O	2:L:123:GLU:C	2.64	0.41
3:H:3:LYS:HE2	3:H:5:VAL:HG11	2.03	0.41
3:H:11:LEU:HB2	3:H:147:PRO:HG3	2.03	0.41
3:H:104:GLY:O	3:H:105:VAL:C	2.64	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
1:E:7:SER:C	1:E:9:ARG:N	2.79	0.41
2:L:34:HIS:HD2	2:L:49:TYR:HA	1.86	0.41
2:L:163:TRP:CZ2	2:L:175:MET:HE3	2.56	0.41
3:H:126:PRO:HD3	3:H:138:LEU:HG	2.03	0.41
2:L:34:HIS:CD2	2:L:49:TYR:HA	2.56	0.40
3:H:138:LEU:HB3	3:H:139:GLY:H	1.72	0.40
3:H:151:THR:OG1	3:H:198:ALA:HB3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:164:ARG:N	3:H:180:LEU:O	2.42	0.40
3:H:177:LEU:C	3:H:177:LEU:CD1	2.94	0.40
2:L:6:ARG:NE	2:L:88:CYS:SG	2.95	0.40
3:H:162:GLY:O	3:H:181:VAL:HA	2.21	0.40
1:E:69:SER:O	1:E:82:ASN:ND2	2.54	0.40
1:E:218:PHE:C	1:E:220:ASP:H	2.29	0.40
3:H:39:GLN:O	3:H:88:ALA:HB1	2.21	0.40
3:H:42:GLY:O	3:H:43:LYS:HB2	2.21	0.40
3:H:188:TRP:HA	3:H:190:SER:N	2.36	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
3:H:34:MET:HE3	3:H:34:MET:HB3	1.86	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	E	269/402 (67%)	207 (77%)	51 (19%)	11 (4%)	2	13
2	L	205/207 (99%)	164 (80%)	36 (18%)	5 (2%)	4	24
3	H	208/221 (94%)	174 (84%)	31 (15%)	3 (1%)	9	36
All	All	682/830 (82%)	545 (80%)	118 (17%)	19 (3%)	4	21

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

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Mol	Chain	Res	Type
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 ⓘ

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%)	206 (91%)	20 (9%)	9	35
2	L	182/182 (100%)	169 (93%)	13 (7%)	13	43
3	H	180/187 (96%)	169 (94%)	11 (6%)	17	49
All	All	588/701 (84%)	544 (92%)	44 (8%)	12	41

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	13	GLU
1	E	15	VAL
1	E	37	ASP
1	E	41	ILE
1	E	58	SER
1	E	70	THR
1	E	96	VAL
1	E	103	ASN

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Mol	Chain	Res	Type
1	E	113	ILE
1	E	115	THR
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	279	VAL
1	E	295	LEU
2	L	22	THR
2	L	30	VAL
2	L	46	ILE
2	L	69	THR
2	L	73	LEU
2	L	76	SER
2	L	92	SER
2	L	97	THR
2	L	123	GLU
2	L	134	CYS
2	L	151	ASP
2	L	189	HIS
2	L	193	THR
3	H	4	LEU
3	H	20	LEU
3	H	24	THR
3	H	38	ARG
3	H	69	ILE
3	H	108	THR
3	H	143	LYS
3	H	159	LEU
3	H	177	LEU
3	H	197	VAL
3	H	206	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	8	ASN
1	E	246	HIS

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Mol	Chain	Res	Type
1	E	277	ASN
2	L	34	HIS
2	L	38	GLN
2	L	137	ASN
2	L	198	HIS
3	H	39	GLN
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 oligosaccharides 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%)	-0.07	3 (1%) 78 57	70, 130, 191, 236	0
2	L	207/207 (100%)	0.03	1 (0%) 87 72	66, 134, 229, 275	0
3	H	212/221 (95%)	0.14	2 (0%) 81 61	67, 148, 276, 316	0
All	All	692/830 (83%)	0.02	6 (0%) 81 61	66, 135, 244, 316	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	H	138	LEU	4.0
1	E	11	PHE	3.6
1	E	247	ALA	2.5
2	L	120	PRO	2.3
3	H	211	SER	2.3
1	E	27	GLY	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 oligosaccharides 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.