



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:17 AM UTC

PDB ID : 1JNL / pdb_00001jnl
Title : Crystal Structure of Fab-Estradiol Complexes
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Deposited on : 2001-07-24
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

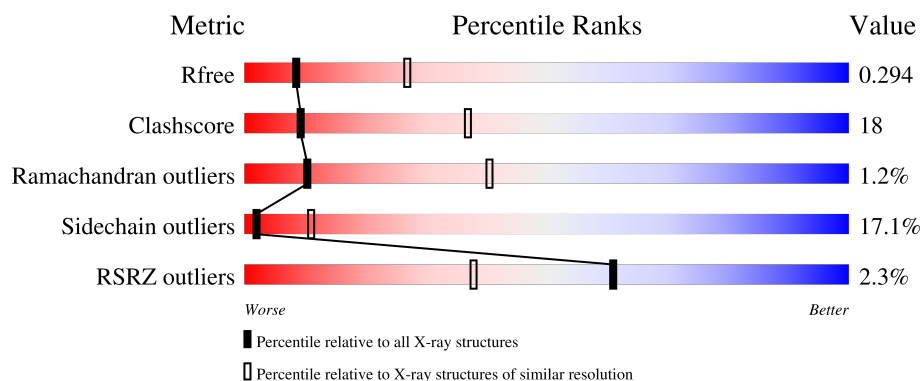
1 Overall quality at a glance

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	L	211	
2	H	216	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3259 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 monoclonal anti-estradiol 17E12E5 immunoglobulin kappa chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	211	Total	C	N	O	S	0	0	0
			1646	1035	277	328	6			

- Molecule 2 is a protein called monoclonal anti-estradiol 17E12E5 immunoglobulin gamma-1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	216	Total	C	N	O	S	0	0	0
			1613	1015	267	324	7			

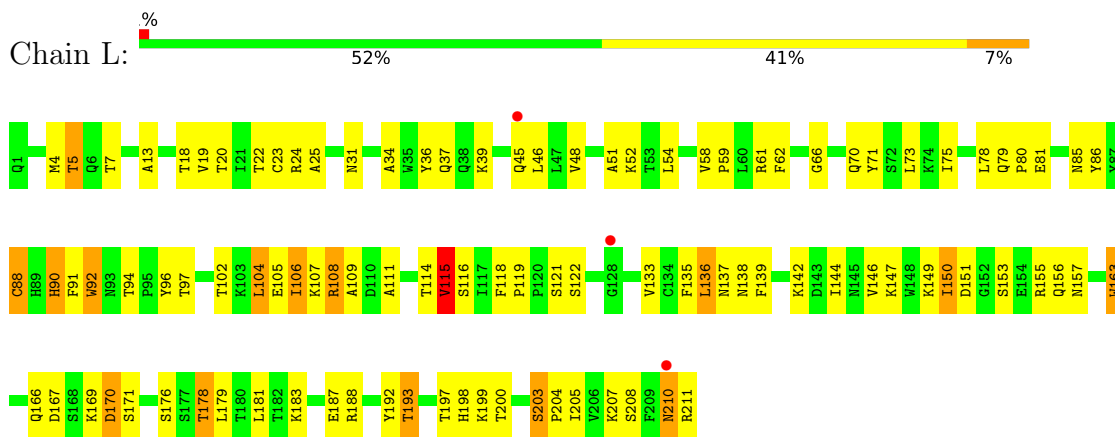
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	222	LYS	-	see remark 999	GB 12658313
H	223	LYS	-	see remark 999	GB 12658313
H	224	ILE	-	see remark 999	GB 12658313
H	227	VAL	-	see remark 999	GB 12658313
H	228	PRO	-	see remark 999	GB 12658313
H	229	ARG	-	see remark 999	GB 12658313
H	230	ASP	-	see remark 999	GB 12658313
H	231	CYS	-	see remark 999	GB 12658313

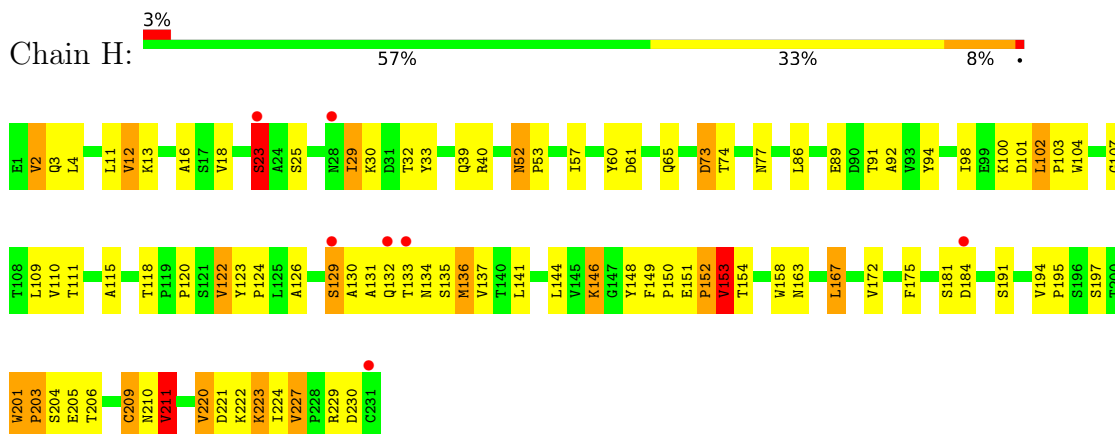
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: monoclonal anti-estradiol 17E12E5 immunoglobulin kappa chain



- Molecule 2: monoclonal anti-estradiol 17E12E5 immunoglobulin gamma-1 chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	70.95Å 143.24Å 53.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00 20.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-3.00) 91.2 (20.00-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.68 (at 2.98Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.195 , 0.290 0.200 , 0.294	Depositor DCC
R_{free} test set	568 reflections (5.23%)	wwPDB-VP
Wilson B-factor (Å ²)	43.7	Xtriage
Anisotropy	0.406	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 67.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3259	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.89% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	L	0.68	0/1687	1.13	14/2293 (0.6%)
2	H	0.66	0/1651	1.18	13/2261 (0.6%)
All	All	0.67	0/3338	1.16	27/4554 (0.6%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	138	ASN	N-CA-C	11.92	128.20	111.52
2	H	61	ASP	N-CA-C	-8.96	97.95	110.31
1	L	203	SER	CA-C-N	7.61	127.67	119.90
1	L	203	SER	C-N-CA	7.61	127.67	119.90
1	L	193	THR	N-CA-C	7.28	120.28	108.34
2	H	52	ASN	CA-C-N	7.27	127.28	119.28
2	H	52	ASN	C-N-CA	7.27	127.28	119.28
1	L	137	ASN	N-CA-C	7.23	120.97	110.28
2	H	204	SER	N-CA-C	-6.76	105.19	113.50
1	L	88	CYS	N-CA-C	-6.69	98.78	109.76
2	H	175	PHE	CA-C-N	6.66	126.62	119.76
2	H	175	PHE	C-N-CA	6.66	126.62	119.76
2	H	211	VAL	N-CA-C	6.45	117.40	108.12
2	H	146	LYS	N-CA-C	6.44	119.10	108.99
1	L	115	VAL	N-CA-C	6.42	118.92	108.97
1	L	92	TRP	N-CA-C	-6.34	103.67	111.40
1	L	79	GLN	CA-C-N	6.31	126.78	119.47
1	L	79	GLN	C-N-CA	6.31	126.78	119.47
2	H	201	TRP	N-CA-C	6.18	120.58	107.91
2	H	13	LYS	N-CA-C	-5.97	101.61	110.20
2	H	153	VAL	N-CA-C	-5.93	100.87	109.29
1	L	170	ASP	N-CA-C	5.79	120.50	113.38
2	H	23	SER	N-CA-C	5.52	119.23	109.96
1	L	97	THR	N-CA-C	5.32	118.39	110.52
1	L	52	LYS	N-CA-C	5.20	120.63	113.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	178	THR	N-CA-C	5.02	116.95	108.02
2	H	129	SER	N-CA-C	5.00	121.45	110.80

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	L	1646	0	1578	66	0
2	H	1613	0	1576	52	0
All	All	3259	0	3154	116	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 18.

All (116) 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:H:206:THR:HG23	2:H:223:LYS:HE3	1.25	1.18
2:H:211:VAL:HG13	2:H:220:VAL:HG13	1.54	0.88
1:L:119:PRO:HG2	2:H:229:ARG:NH2	1.95	0.81
1:L:150:ILE:HG23	1:L:151:ASP:H	1.48	0.78
2:H:122:VAL:HG22	2:H:222:LYS:HG3	1.67	0.76
2:H:141:LEU:HD12	2:H:224:ILE:HG21	1.68	0.76
2:H:30:LYS:HB3	2:H:74:THR:HG21	1.69	0.74
1:L:155:ARG:NH1	1:L:157:ASN:HB3	2.05	0.70
1:L:108:ARG:HG2	1:L:109:ALA:N	2.06	0.69
1:L:193:THR:HG23	1:L:208:SER:HB3	1.73	0.69
1:L:115:VAL:HB	1:L:136:LEU:HD13	1.75	0.69
2:H:206:THR:CG2	2:H:223:LYS:HE3	2.15	0.68
2:H:12:VAL:HG21	2:H:86:LEU:HD13	1.77	0.67
1:L:166:GLN:NE2	1:L:171:SER:HB3	2.10	0.67
2:H:102:LEU:HG	2:H:103:PRO:HD2	1.76	0.66
2:H:100:LYS:HG2	2:H:101:ASP:H	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:198:HIS:CG	1:L:199:LYS:H	2.18	0.61
1:L:108:ARG:HH21	1:L:111:ALA:HB2	1.65	0.60
1:L:90:HIS:HD2	1:L:92:TRP:H	1.48	0.60
1:L:19:VAL:HG22	1:L:75:ILE:HB	1.84	0.60
1:L:66:GLY:HA3	1:L:71:TYR:HA	1.83	0.60
1:L:90:HIS:C	1:L:90:HIS:CD2	2.80	0.59
1:L:142:LYS:O	1:L:144:ILE:HG22	2.04	0.58
1:L:19:VAL:CG1	1:L:78:LEU:HD11	2.34	0.57
1:L:183:LYS:O	1:L:187:GLU:HG2	2.05	0.57
2:H:136:MET:HG3	2:H:195:PRO:HA	1.87	0.57
1:L:163:TRP:N	1:L:163:TRP:CD1	2.72	0.57
1:L:139:PHE:CZ	1:L:144:ILE:HG21	2.41	0.56
2:H:4:LEU:HD12	2:H:4:LEU:N	2.21	0.56
1:L:80:PRO:HA	1:L:106:ILE:HD12	1.88	0.56
2:H:4:LEU:HA	2:H:23:SER:O	2.05	0.56
1:L:4:MET:HE1	1:L:25:ALA:HB2	1.87	0.56
2:H:94:TYR:O	2:H:107:GLY:HA2	2.06	0.55
2:H:33:TYR:CE2	2:H:52:ASN:HB2	2.41	0.55
2:H:2:VAL:HA	2:H:25:SER:O	2.07	0.55
2:H:126:ALA:O	2:H:229:ARG:NH2	2.40	0.54
2:H:39:GLN:O	2:H:92:ALA:HB1	2.08	0.54
1:L:59:PRO:HG2	1:L:62:PHE:CD2	2.43	0.54
1:L:37:GLN:HA	1:L:85:ASN:O	2.08	0.53
1:L:24:ARG:HD3	1:L:70:GLN:HE21	1.73	0.53
2:H:120:PRO:HB3	2:H:148:TYR:HB3	1.91	0.52
2:H:122:VAL:HG11	2:H:211:VAL:CG1	2.39	0.52
2:H:172:VAL:HA	2:H:191:SER:O	2.09	0.52
1:L:167:ASP:OD1	1:L:169:LYS:HD2	2.08	0.52
2:H:151:GLU:OE1	2:H:152:PRO:HA	2.09	0.52
2:H:4:LEU:HD12	2:H:4:LEU:H	1.73	0.52
1:L:150:ILE:HG23	1:L:151:ASP:N	2.21	0.52
1:L:166:GLN:HG2	1:L:171:SER:O	2.10	0.52
1:L:167:ASP:HB3	1:L:170:ASP:OD1	2.10	0.52
2:H:154:THR:O	2:H:211:VAL:HA	2.09	0.52
1:L:210:ASN:O	1:L:211:ARG:HB3	2.09	0.51
1:L:59:PRO:HG2	1:L:62:PHE:HD2	1.74	0.51
2:H:223:LYS:HD3	2:H:227:VAL:HG13	1.93	0.50
2:H:30:LYS:HB3	2:H:74:THR:CG2	2.41	0.50
2:H:122:VAL:HG11	2:H:211:VAL:HG11	1.95	0.49
1:L:85:ASN:HA	1:L:102:THR:O	2.12	0.49
1:L:150:ILE:HG22	1:L:153:SER:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:60:TYR:CD1	2:H:60:TYR:N	2.81	0.48
1:L:205:ILE:HD12	1:L:205:ILE:N	2.28	0.48
2:H:91:THR:HG23	2:H:111:THR:HA	1.95	0.48
2:H:120:PRO:CA	2:H:148:TYR:HB3	2.43	0.48
2:H:91:THR:HA	2:H:110:VAL:O	2.14	0.48
1:L:155:ARG:HH12	1:L:157:ASN:HB3	1.78	0.48
1:L:179:LEU:HG	1:L:181:LEU:CD1	2.44	0.48
2:H:11:LEU:HD22	2:H:150:PRO:HG3	1.95	0.47
2:H:201:TRP:HA	2:H:203:PRO:HA	1.51	0.47
1:L:115:VAL:HG22	1:L:207:LYS:HG3	1.96	0.47
1:L:192:TYR:O	1:L:208:SER:HB2	2.15	0.47
1:L:104:LEU:C	1:L:104:LEU:HD12	2.40	0.46
2:H:73:ASP:OD1	2:H:73:ASP:O	2.34	0.46
2:H:122:VAL:HG13	2:H:220:VAL:HG21	1.96	0.46
2:H:130:ALA:O	2:H:132:GLN:N	2.49	0.46
1:L:13:ALA:O	1:L:107:LYS:N	2.45	0.46
1:L:36:TYR:O	1:L:86:TYR:HA	2.16	0.46
1:L:20:THR:HA	1:L:73:LEU:O	2.16	0.45
1:L:88:CYS:O	1:L:88:CYS:SG	2.73	0.45
2:H:223:LYS:O	2:H:227:VAL:HG22	2.17	0.45
1:L:136:LEU:N	1:L:136:LEU:HD22	2.31	0.45
1:L:166:GLN:HG2	1:L:171:SER:C	2.41	0.45
1:L:139:PHE:HE1	1:L:142:LYS:HA	1.82	0.45
1:L:198:HIS:CG	1:L:199:LYS:N	2.85	0.45
1:L:90:HIS:CD2	1:L:92:TRP:H	2.32	0.45
1:L:34:ALA:HB1	1:L:46:LEU:HD11	1.99	0.45
1:L:91:PHE:HA	1:L:96:TYR:CD1	2.52	0.44
2:H:104:TRP:N	2:H:104:TRP:CD1	2.84	0.44
2:H:16:ALA:O	2:H:86:LEU:HG	2.17	0.44
1:L:121:SER:OG	2:H:123:TYR:HB3	2.17	0.44
1:L:90:HIS:CD2	1:L:91:PHE:N	2.86	0.44
1:L:115:VAL:HG13	1:L:205:ILE:HG21	2.00	0.44
1:L:105:GLU:HG2	1:L:106:ILE:N	2.32	0.44
2:H:181:SER:O	2:H:184:ASP:HB2	2.17	0.44
2:H:29:ILE:CG2	2:H:53:PRO:HG3	2.49	0.43
1:L:19:VAL:HG13	1:L:78:LEU:HD11	1.99	0.43
1:L:133:VAL:HG22	1:L:178:THR:OG1	2.18	0.43
1:L:108:ARG:NH2	1:L:111:ALA:HB2	2.33	0.43
1:L:19:VAL:HG11	1:L:78:LEU:HD11	2.00	0.43
1:L:59:PRO:HB2	1:L:61:ARG:HG2	2.01	0.43
2:H:12:VAL:HG21	2:H:86:LEU:CD1	2.47	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:4:MET:SD	1:L:90:HIS:HB3	2.59	0.42
1:L:5:THR:O	1:L:5:THR:HG22	2.19	0.42
2:H:137:VAL:HG12	2:H:194:VAL:O	2.19	0.42
2:H:158:TRP:CZ3	2:H:209:CYS:HB2	2.54	0.42
2:H:153:VAL:HG13	2:H:154:THR:N	2.33	0.42
2:H:4:LEU:HD21	2:H:98:ILE:HG23	2.01	0.41
2:H:124:PRO:HB2	2:H:224:ILE:HD13	2.01	0.41
1:L:31:ASN:HD22	1:L:51:ALA:HB3	1.85	0.41
1:L:39:LYS:HE3	1:L:39:LYS:HB2	1.75	0.41
1:L:118:PHE:CE2	1:L:135:PHE:CD2	3.08	0.41
1:L:54:LEU:HD11	1:L:58:VAL:HB	2.02	0.41
2:H:120:PRO:CB	2:H:148:TYR:HB3	2.50	0.41
2:H:122:VAL:O	2:H:222:LYS:HE2	2.21	0.41
1:L:166:GLN:HE21	1:L:171:SER:HB3	1.84	0.40
1:L:203:SER:HA	1:L:204:PRO:HD2	1.86	0.40
2:H:115:ALA:HB3	2:H:149:PHE:CE1	2.55	0.40
2:H:163:ASN:ND2	2:H:167:LEU:HD12	2.36	0.40
1:L:118:PHE:HE2	1:L:135:PHE:CD2	2.39	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	L	209/211 (99%)	199 (95%)	9 (4%)	1 (0%)	24	60
2	H	214/216 (99%)	196 (92%)	14 (6%)	4 (2%)	6	30
All	All	423/427 (99%)	395 (93%)	23 (5%)	5 (1%)	10	40

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	150	ILE
2	H	129	SER
2	H	131	ALA
2	H	135	SER
2	H	203	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	185/186 (100%)	157 (85%)	28 (15%)	3	14
2	H	183/186 (98%)	148 (81%)	35 (19%)	1	8
All	All	368/372 (99%)	305 (83%)	63 (17%)	2	11

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	5	THR
1	L	7	THR
1	L	18	THR
1	L	22	THR
1	L	23	CYS
1	L	45	GLN
1	L	48	VAL
1	L	81	GLU
1	L	90	HIS
1	L	94	THR
1	L	104	LEU
1	L	106	ILE
1	L	108	ARG
1	L	114	THR
1	L	115	VAL
1	L	116	SER
1	L	122	SER
1	L	136	LEU
1	L	146	VAL

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Mol	Chain	Res	Type
1	L	147	LYS
1	L	149	LYS
1	L	156	GLN
1	L	163	TRP
1	L	176	SER
1	L	188	ARG
1	L	197	THR
1	L	200	THR
1	L	210	ASN
2	H	2	VAL
2	H	3	GLN
2	H	12	VAL
2	H	18	VAL
2	H	23	SER
2	H	29	ILE
2	H	32	THR
2	H	40	ARG
2	H	57	ILE
2	H	65	GLN
2	H	73	ASP
2	H	77	ASN
2	H	89	GLU
2	H	102	LEU
2	H	109	LEU
2	H	118	THR
2	H	122	VAL
2	H	133	THR
2	H	134	ASN
2	H	136	MET
2	H	144	LEU
2	H	146	LYS
2	H	152	PRO
2	H	153	VAL
2	H	167	LEU
2	H	197	SER
2	H	205	GLU
2	H	209	CYS
2	H	210	ASN
2	H	211	VAL
2	H	220	VAL
2	H	221	ASP
2	H	223	LYS

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Mol	Chain	Res	Type
2	H	227	VAL
2	H	230	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	31	ASN
1	L	40	GLN
1	L	70	GLN
1	L	85	ASN
1	L	90	HIS
1	L	156	GLN
2	H	65	GLN
2	H	82	GLN
2	H	173	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 ⓘ

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	L	211/211 (100%)	-0.00	3 (1%) 73 51	10, 20, 68, 90	0
2	H	216/216 (100%)	-0.07	7 (3%) 50 29	10, 20, 79, 90	0
All	All	427/427 (100%)	-0.04	10 (2%) 61 38	10, 20, 76, 90	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	133	THR	5.1
2	H	184	ASP	4.9
2	H	129	SER	4.9
2	H	132	GLN	3.1
2	H	231	CYS	2.7
1	L	128	GLY	2.5
1	L	210	ASN	2.4
1	L	45	GLN	2.1
2	H	23	SER	2.0
2	H	28	ASN	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.