



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:47 AM UTC

PDB ID : 6WMW / pdb_00006wmw
Title : GFRAL receptor bound with two antibody Fabs (3P10, 25M22)
Authors : White, A.; Lakshminarasimhan, D.; Olland, A.; Suto, R.K.
Deposited on : 2020-04-21
Resolution : 2.91 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

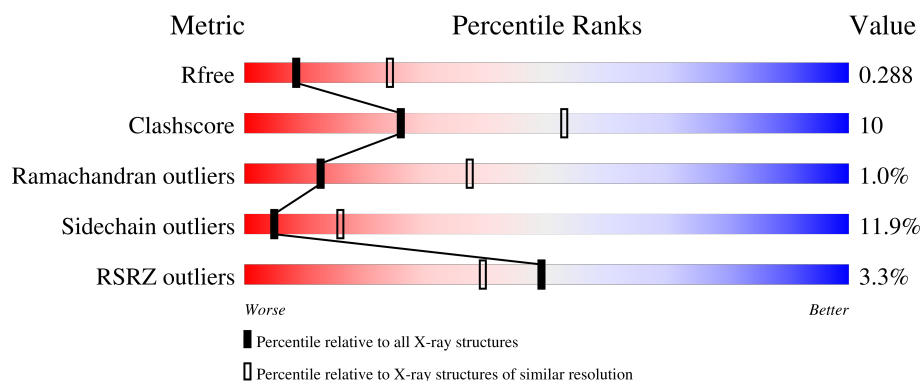
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.91 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	2995 (2.94-2.90)
Clashscore	190562	3213 (2.94-2.90)
Ramachandran outliers	187476	3128 (2.94-2.90)
Sidechain outliers	187428	3130 (2.94-2.90)
RSRZ outliers	180081	2995 (2.94-2.90)

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	B	245	
2	H	225	
3	L	218	
4	M	232	
5	N	241	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8097 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 GDNF family receptor alpha-like.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	192	Total	C	N	O	S	0	0	0
			1496	921	266	284	25			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	107	HIS	-	expression tag	UNP Q6UXV0
B	108	HIS	-	expression tag	UNP Q6UXV0
B	109	HIS	-	expression tag	UNP Q6UXV0
B	110	HIS	-	expression tag	UNP Q6UXV0
B	111	HIS	-	expression tag	UNP Q6UXV0
B	112	HIS	-	expression tag	UNP Q6UXV0
B	113	GLY	-	expression tag	UNP Q6UXV0
B	114	GLY	-	expression tag	UNP Q6UXV0

- Molecule 2 is a protein called FAB3P10 heavy chain fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	212	Total	C	N	O	S	0	0	0
			1592	1015	258	314	5			

- Molecule 3 is a protein called FAB3P10 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	218	Total	C	N	O	S	0	0	0
			1660	1040	278	334	8			

- Molecule 4 is a protein called FAB25M22 heavy chain fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	M	222	Total	C	N	O	S	0	0	0
			1671	1060	279	325	7			

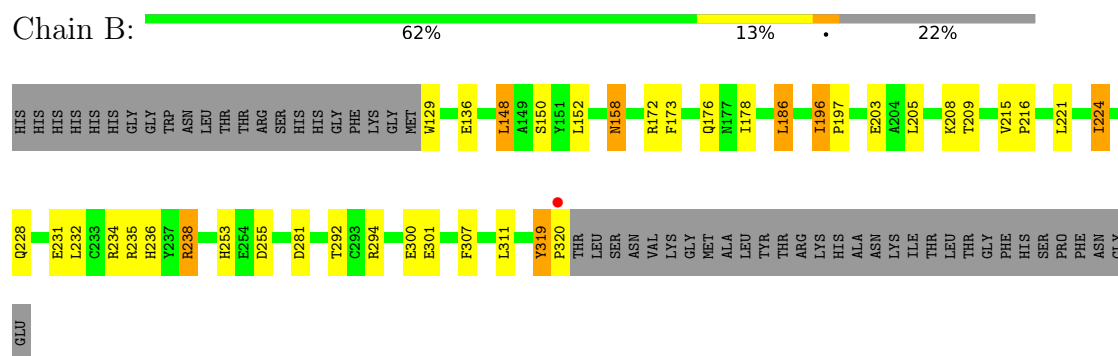
- Molecule 5 is a protein called FAB25M22 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	N	216	Total	C	N	O	S	0	0	0
			1678	1062	275	337	4			

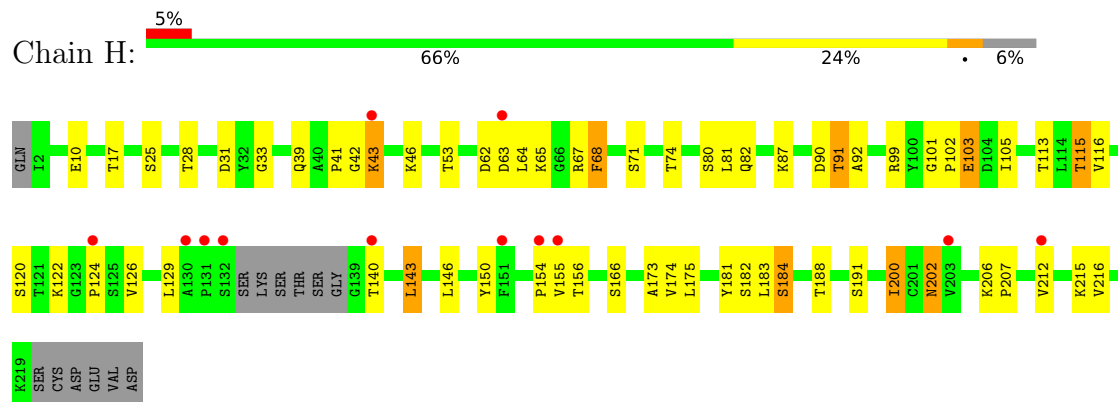
3 Residue-property plots [i](#)

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.

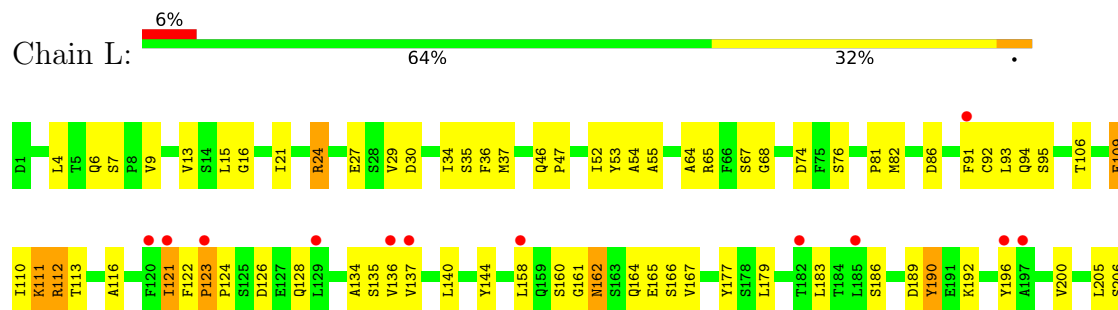
- Molecule 1: GDNF family receptor alpha-like



- Molecule 2: FAB3P10 heavy chain fragment

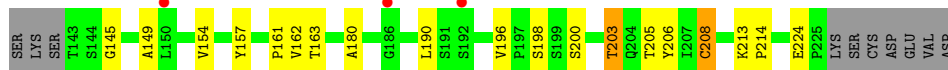
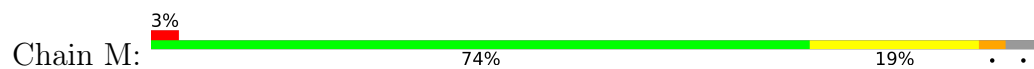


- Molecule 3: FAB3P10 light chain

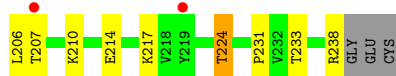




- Molecule 4: FAB25M22 heavy chain fragment



- Molecule 5: FAB25M22 light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.50Å 116.05Å 227.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.70 – 2.91 47.70 – 2.91	Depositor EDS
% Data completeness (in resolution range)	99.4 (47.70-2.91) 99.6 (47.70-2.91)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.8.0230	Depositor
R, R_{free}	0.222 , 0.290 0.224 , 0.288	Depositor DCC
R_{free} test set	1591 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	64.6	Xtriage
Anisotropy	0.091	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 27.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8097	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.62% 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	B	0.72	0/1523	1.06	0/2060
2	H	0.74	0/1633	0.99	0/2230
3	L	0.78	0/1698	1.03	2/2303 (0.1%)
4	M	0.68	0/1713	0.93	1/2331 (0.0%)
5	N	0.69	0/1715	0.96	1/2329 (0.0%)
All	All	0.72	0/8282	0.99	4/11253 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	L	0	1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	M	58	THR	CA-CB-OG1	-5.92	100.72	109.60
3	L	64	ALA	CA-C-N	5.49	128.18	120.28
3	L	64	ALA	C-N-CA	5.49	128.18	120.28
5	N	140	PRO	CB-CA-C	-5.02	106.83	111.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	L	112	ARG	Sidechain

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	B	1496	0	1434	23	0
2	H	1592	0	1548	35	0
3	L	1660	0	1596	42	0
4	M	1671	0	1642	26	0
5	N	1678	0	1633	35	0
All	All	8097	0	7853	152	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (152) 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)
4:M:104:LEU:O	4:M:105:ARG:HD3	1.69	0.92
4:M:87:THR:O	4:M:123:VAL:HG21	1.77	0.83
2:H:67:ARG:NH1	2:H:90:ASP:OD2	2.13	0.81
5:N:214:GLU:HG2	5:N:238:ARG:HH11	1.50	0.76
5:N:214:GLU:HG2	5:N:238:ARG:NH1	2.01	0.76
1:B:232:LEU:O	1:B:236:HIS:ND1	2.18	0.76
2:H:87:LYS:O	2:H:116:VAL:HG11	1.86	0.75
5:N:135:ARG:NH1	5:N:197:ASP:O	2.20	0.73
3:L:30:ASP:OD1	3:L:35:SER:OG	2.04	0.72
2:H:124:PRO:HB3	2:H:150:TYR:HB3	1.72	0.70
3:L:21:ILE:HD12	3:L:106:THR:HG21	1.74	0.69
1:B:224:ILE:HD12	1:B:311:LEU:HD11	1.75	0.69
2:H:91:THR:HG23	2:H:115:THR:HA	1.77	0.67
2:H:129:LEU:HD23	3:L:122:PHE:CG	2.30	0.66
5:N:94:SER:O	5:N:96:THR:N	2.29	0.65
1:B:319:TYR:HB3	1:B:320:PRO:CD	2.27	0.64
2:H:143:LEU:HD13	2:H:216:VAL:HG11	1.79	0.64
3:L:121:ILE:HG21	3:L:211:LYS:O	1.97	0.64
5:N:88:ARG:NH1	5:N:109:ASP:OD2	2.30	0.64
3:L:65:ARG:NH2	3:L:86:ASP:OD1	2.29	0.64
5:N:33:LEU:HD21	5:N:41:ALA:HB1	1.80	0.64
1:B:224:ILE:HD13	1:B:307:PHE:CD1	2.34	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:M:208:CYS:O	4:M:208:CYS:SG	2.56	0.63
4:M:98:ARG:O	4:M:112:MET:HA	2.00	0.62
5:N:88:ARG:HH12	5:N:109:ASP:CG	2.08	0.61
1:B:203:GLU:OE2	4:M:106:ARG:HD3	2.01	0.60
2:H:99:ARG:O	2:H:101:GLY:HA2	2.01	0.60
1:B:319:TYR:CB	1:B:320:PRO:CD	2.79	0.60
2:H:183:LEU:C	2:H:183:LEU:HD12	2.27	0.59
4:M:131:PRO:HB3	4:M:157:TYR:HB3	1.85	0.58
3:L:123:PRO:HB3	3:L:213:PHE:CE2	2.38	0.58
1:B:221:LEU:HD11	1:B:300:GLU:HB3	1.86	0.58
3:L:165:GLU:HB2	3:L:179:LEU:HD11	1.86	0.57
5:N:60:LEU:HD22	5:N:98:PHE:CG	2.39	0.57
4:M:104:LEU:O	4:M:105:ARG:CD	2.50	0.56
2:H:43:LYS:O	2:H:43:LYS:HG3	2.06	0.56
1:B:319:TYR:CB	1:B:320:PRO:HD2	2.36	0.56
4:M:12:VAL:O	4:M:123:VAL:HA	2.07	0.55
2:H:200:ILE:HG13	2:H:215:LYS:HA	1.89	0.55
1:B:234:ARG:NH1	1:B:238:ARG:HH21	2.03	0.55
3:L:121:ILE:HG13	3:L:122:PHE:N	2.21	0.55
5:N:145:PHE:O	5:N:159:VAL:HG22	2.06	0.55
4:M:36:TRP:CD2	4:M:81:MET:HE3	2.42	0.55
2:H:64:LEU:HG	2:H:68:PHE:CE1	2.43	0.54
2:H:33:GLY:HA2	2:H:53:THR:HG23	1.88	0.54
4:M:36:TRP:CG	4:M:81:MET:HE3	2.42	0.54
2:H:42:GLY:O	2:H:43:LYS:HG2	2.07	0.54
2:H:129:LEU:HD23	3:L:122:PHE:CD2	2.42	0.54
5:N:159:VAL:HG12	5:N:206:LEU:HB3	1.90	0.54
3:L:123:PRO:HB3	3:L:213:PHE:CZ	2.44	0.53
3:L:162:ASN:OD1	3:L:162:ASN:N	2.41	0.53
3:L:140:LEU:HD21	3:L:200:VAL:HG13	1.91	0.53
4:M:81:MET:SD	4:M:83:LEU:HD21	2.49	0.53
1:B:231:GLU:O	1:B:235:ARG:HG3	2.08	0.53
3:L:16:GLY:HA2	3:L:81:PRO:HB2	1.91	0.53
3:L:29:VAL:HG12	3:L:29:VAL:O	2.10	0.52
5:N:53:ASN:HB3	5:N:55:ASP:OD1	2.10	0.52
1:B:186:LEU:HD21	1:B:205:LEU:HD22	1.91	0.52
3:L:37:MET:SD	3:L:92:CYS:HB2	2.49	0.52
3:L:82:MET:HE3	3:L:110:ILE:CD1	2.40	0.51
3:L:65:ARG:HD3	3:L:81:PRO:O	2.10	0.51
4:M:180:ALA:HA	4:M:190:LEU:HB3	1.93	0.51
1:B:172:ARG:CD	1:B:232:LEU:HD21	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:206:LYS:N	2:H:207:PRO:CD	2.74	0.51
4:M:112:MET:HG3	5:N:63:TYR:OH	2.10	0.50
4:M:149:ALA:HB3	5:N:143:PHE:CE2	2.46	0.50
1:B:196:ILE:N	1:B:197:PRO:HD2	2.26	0.50
3:L:190:TYR:HD1	3:L:196:TYR:HH	1.58	0.50
4:M:213:LYS:N	4:M:214:PRO:CD	2.74	0.50
3:L:82:MET:HE3	3:L:110:ILE:HD13	1.92	0.50
2:H:71:SER:OG	2:H:80:SER:HB2	2.11	0.50
5:N:192:GLU:O	5:N:193:GLN:C	2.54	0.49
5:N:118:SER:HA	5:N:123:TYR:CD1	2.47	0.49
5:N:147:PRO:HD3	5:N:159:VAL:HG23	1.95	0.48
2:H:129:LEU:HB3	3:L:122:PHE:CD2	2.49	0.48
5:N:155:GLY:O	5:N:210:LYS:HB2	2.14	0.48
5:N:62:TRP:HB2	5:N:75:ILE:HB	1.96	0.48
3:L:136:VAL:N	3:L:183:LEU:O	2.45	0.47
5:N:163:LEU:HD12	5:N:163:LEU:N	2.29	0.47
5:N:66:LYS:HD3	5:N:108:GLU:O	2.15	0.47
1:B:292:THR:HB	1:B:294:ARG:H	1.80	0.47
3:L:186:SER:OG	3:L:189:ASP:HB2	2.14	0.47
4:M:88:SER:HA	4:M:123:VAL:CG2	2.45	0.47
3:L:121:ILE:HD11	3:L:213:PHE:CD1	2.50	0.47
1:B:319:TYR:HB3	1:B:320:PRO:HD3	1.96	0.46
3:L:4:LEU:HA	3:L:24:ARG:O	2.15	0.46
3:L:196:TYR:HB2	3:L:213:PHE:CE1	2.51	0.46
1:B:172:ARG:HD3	1:B:232:LEU:HD21	1.97	0.46
5:N:64:LEU:HD12	5:N:65:GLN:N	2.31	0.46
5:N:161:CYS:HB2	5:N:175:TRP:CH2	2.51	0.46
2:H:173:ALA:HA	2:H:183:LEU:HB3	1.98	0.46
3:L:36:PHE:O	3:L:94:GLN:HA	2.16	0.46
1:B:319:TYR:HB3	1:B:320:PRO:HD2	1.95	0.45
2:H:39:GLN:O	2:H:92:ALA:HB1	2.16	0.45
3:L:55:ALA:O	3:L:68:GLY:HA3	2.15	0.45
2:H:202:ASN:H	2:H:202:ASN:HD22	1.65	0.45
1:B:292:THR:HG22	3:L:34:ILE:HD13	1.99	0.45
5:N:146:PRO:HA	5:N:159:VAL:HG23	1.98	0.45
2:H:41:PRO:C	2:H:43:LYS:H	2.23	0.45
4:M:47:TRP:CZ2	4:M:49:GLY:HA2	2.52	0.45
3:L:128:GLN:CB	3:L:134:ALA:HB2	2.46	0.45
5:N:135:ARG:NH1	5:N:199:THR:HG23	2.32	0.44
5:N:163:LEU:HD13	5:N:202:LEU:HB3	1.98	0.44
2:H:102:PRO:HB2	2:H:103:GLU:HG2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:212:VAL:O	2:H:212:VAL:HG13	2.18	0.44
4:M:127:SER:O	4:M:128:THR:C	2.61	0.44
3:L:196:TYR:CB	3:L:213:PHE:CE1	3.01	0.44
4:M:196:VAL:HG11	4:M:206:TYR:CZ	2.53	0.44
2:H:67:ARG:HH12	2:H:90:ASP:CG	2.26	0.44
2:H:175:LEU:HD13	2:H:181:TYR:CE1	2.53	0.44
4:M:190:LEU:N	4:M:190:LEU:HD23	2.33	0.44
5:N:36:ASN:OD1	5:N:134:LYS:HB3	2.18	0.44
5:N:28:GLN:HA	5:N:44:SER:O	2.18	0.43
5:N:214:GLU:CG	5:N:238:ARG:HH11	2.25	0.43
5:N:224:THR:HG23	5:N:231:PRO:HG3	2.00	0.43
1:B:301:GLU:HG3	2:H:102:PRO:HG3	1.99	0.43
3:L:167:VAL:HG22	3:L:179:LEU:HD13	2.00	0.43
5:N:161:CYS:HB2	5:N:175:TRP:CZ2	2.53	0.43
2:H:129:LEU:HD21	3:L:137:VAL:HG21	2.01	0.43
1:B:224:ILE:O	1:B:228:GLN:HG2	2.19	0.43
2:H:202:ASN:HD22	2:H:202:ASN:N	2.16	0.43
5:N:116:PHE:HB2	5:N:125:PHE:CD1	2.54	0.43
3:L:116:ALA:HB1	3:L:205:LEU:HG	2.01	0.43
4:M:105:ARG:O	4:M:106:ARG:C	2.62	0.43
4:M:200:SER:HA	4:M:203:THR:HG23	2.01	0.43
3:L:109:GLU:OE1	3:L:177:TYR:OH	2.34	0.42
3:L:111:LYS:HA	3:L:144:TYR:OH	2.19	0.42
5:N:202:LEU:HD23	5:N:203:SER:N	2.34	0.42
3:L:15:LEU:HD12	3:L:15:LEU:H	1.84	0.42
2:H:183:LEU:HD12	2:H:184:SER:N	2.34	0.42
1:B:148:LEU:HD22	1:B:152:LEU:CD1	2.49	0.42
2:H:174:VAL:HG22	2:H:182:SER:O	2.20	0.42
5:N:50:SER:HA	5:N:95:GLY:O	2.19	0.42
3:L:112:ARG:HH11	3:L:112:ARG:HG2	1.85	0.42
4:M:12:VAL:HG23	4:M:123:VAL:HG12	2.02	0.42
4:M:47:TRP:CH2	4:M:49:GLY:HA2	2.55	0.42
4:M:14:PRO:HA	4:M:15:GLY:HA2	1.75	0.42
2:H:202:ASN:N	2:H:202:ASN:ND2	2.68	0.42
5:N:43:ILE:HG12	5:N:129:THR:HG21	2.03	0.41
3:L:6:GLN:HE22	3:L:91:PHE:HA	1.85	0.41
3:L:53:TYR:O	3:L:54:ALA:C	2.63	0.41
3:L:124:PRO:HD3	3:L:136:VAL:HG22	2.03	0.41
3:L:46:GLN:HB3	3:L:47:PRO:HD2	2.02	0.41
2:H:63:ASP:N	2:H:63:ASP:OD1	2.53	0.41
1:B:173:PHE:CD1	1:B:173:PHE:C	2.98	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:215:VAL:HA	1:B:216:PRO:HA	1.81	0.41
2:H:124:PRO:CB	2:H:150:TYR:HB3	2.48	0.41
3:L:13:VAL:CG1	3:L:82:MET:HG3	2.51	0.41
2:H:91:THR:CG2	2:H:115:THR:HA	2.49	0.40
5:N:64:LEU:HB2	5:N:74:LEU:HD11	2.03	0.40
2:H:28:THR:HB	2:H:31:ASP:CG	2.47	0.40
4:M:36:TRP:CB	4:M:81:MET:CE	2.99	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	B	190/245 (78%)	177 (93%)	10 (5%)	3 (2%)	7	25
2	H	208/225 (92%)	166 (80%)	42 (20%)	0	100	100
3	L	216/218 (99%)	189 (88%)	24 (11%)	3 (1%)	9	29
4	M	218/232 (94%)	191 (88%)	25 (12%)	2 (1%)	14	40
5	N	214/241 (89%)	192 (90%)	20 (9%)	2 (1%)	14	40
All	All	1046/1161 (90%)	915 (88%)	121 (12%)	10 (1%)	12	37

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	319	TYR
5	N	95	GLY
5	N	141	SER
1	B	158	ASN
3	L	208	PRO
3	L	161	GLY
1	B	255	ASP

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Mol	Chain	Res	Type
4	M	41	PRO
3	L	123	PRO
4	M	145	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	173/217 (80%)	158 (91%)	15 (9%)	9	28
2	H	176/191 (92%)	145 (82%)	31 (18%)	2	6
3	L	187/189 (99%)	162 (87%)	25 (13%)	4	12
4	M	185/196 (94%)	161 (87%)	24 (13%)	4	13
5	N	193/214 (90%)	179 (93%)	14 (7%)	13	37
All	All	914/1007 (91%)	805 (88%)	109 (12%)	5	16

All (109) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	129	TRP
1	B	136	GLU
1	B	148	LEU
1	B	150	SER
1	B	158	ASN
1	B	176	GLN
1	B	178	ILE
1	B	186	LEU
1	B	196	ILE
1	B	208	LYS
1	B	209	THR
1	B	224	ILE
1	B	238	ARG
1	B	253	HIS
1	B	281	ASP
2	H	10	GLU

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Mol	Chain	Res	Type
2	H	17	THR
2	H	25	SER
2	H	43	LYS
2	H	46	LYS
2	H	62	ASP
2	H	65	LYS
2	H	68	PHE
2	H	74	THR
2	H	81	LEU
2	H	82	GLN
2	H	91	THR
2	H	103	GLU
2	H	105	ILE
2	H	113	THR
2	H	115	THR
2	H	120	SER
2	H	122	LYS
2	H	126	VAL
2	H	140	THR
2	H	143	LEU
2	H	146	LEU
2	H	154	PRO
2	H	155	VAL
2	H	156	THR
2	H	166	SER
2	H	184	SER
2	H	188	THR
2	H	191	SER
2	H	200	ILE
2	H	202	ASN
3	L	7	SER
3	L	9	VAL
3	L	24	ARG
3	L	27	GLU
3	L	52	ILE
3	L	67	SER
3	L	74	ASP
3	L	76	SER
3	L	93	LEU
3	L	95	SER
3	L	109	GLU
3	L	111	LYS

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Mol	Chain	Res	Type
3	L	113	THR
3	L	121	ILE
3	L	126	ASP
3	L	135	SER
3	L	158	LEU
3	L	160	SER
3	L	162	ASN
3	L	164	GLN
3	L	166	SER
3	L	190	TYR
3	L	192	LYS
3	L	206	SER
3	L	215	ARG
4	M	10	ASP
4	M	11	LEU
4	M	21	SER
4	M	37	MET
4	M	65	LYS
4	M	71	THR
4	M	77	SER
4	M	87	THR
4	M	88	SER
4	M	107	THR
4	M	112	MET
4	M	120	SER
4	M	125	SER
4	M	127	SER
4	M	139	SER
4	M	154	VAL
4	M	161	PRO
4	M	162	VAL
4	M	163	THR
4	M	198	SER
4	M	203	THR
4	M	205	THR
4	M	208	CYS
4	M	224	GLU
5	N	23	ASP
5	N	27	THR
5	N	29	THR
5	N	40	GLN
5	N	83	SER

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Mol	Chain	Res	Type
5	N	121	LEU
5	N	141	SER
5	N	159	VAL
5	N	170	GLU
5	N	185	ASN
5	N	207	THR
5	N	217	LYS
5	N	224	THR
5	N	233	THR

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

Mol	Chain	Res	Type
2	H	39	GLN
2	H	84	ASN
3	L	42	GLN
4	M	59	ASN
5	N	72	GLN
5	N	226	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers

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

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	B	192/245 (78%)	-0.11	1 (0%) 87 83	35, 59, 94, 137	0
2	H	212/225 (94%)	0.47	12 (5%) 29 23	37, 82, 134, 151	0
3	L	218/218 (100%)	0.06	12 (5%) 30 24	30, 60, 128, 160	0
4	M	222/232 (95%)	0.17	6 (2%) 56 46	39, 74, 124, 143	0
5	N	216/241 (89%)	0.36	4 (1%) 66 58	42, 83, 157, 179	0
All	All	1060/1161 (91%)	0.20	35 (3%) 49 39	30, 71, 135, 179	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	203	VAL	3.9
4	M	150	LEU	3.7
4	M	138	PRO	3.3
2	H	132	SER	3.3
2	H	130	ALA	3.2
1	B	320	PRO	3.0
3	L	121	ILE	2.9
2	H	43	LYS	2.7
3	L	136	VAL	2.7
3	L	137	VAL	2.7
4	M	137	ALA	2.6
2	H	212	VAL	2.6
3	L	120	PHE	2.6
4	M	192	SER	2.6
4	M	135	PRO	2.6
3	L	91	PHE	2.5
2	H	63	ASP	2.5
2	H	124	PRO	2.4
5	N	207	THR	2.3
2	H	131	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
3	L	123	PRO	2.3
3	L	196	TYR	2.3
5	N	149	ASP	2.2
2	H	140	THR	2.2
5	N	219	TYR	2.2
2	H	154	PRO	2.2
3	L	129	LEU	2.2
3	L	182	THR	2.2
2	H	155	VAL	2.1
3	L	185	LEU	2.1
2	H	151	PHE	2.1
3	L	197	ALA	2.1
5	N	177	VAL	2.0
4	M	186	GLY	2.0
3	L	158	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 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.