



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:42 PM UTC

PDB ID : 8ATH / pdb_00008ath
Title : CRYSTAL STRUCTURE OF LAMP1 IN COMPLEX WITH FAB-B.
Authors : Mathieu, M.; Dupuy, A.
Deposited on : 2022-08-23
Resolution : 2.37 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

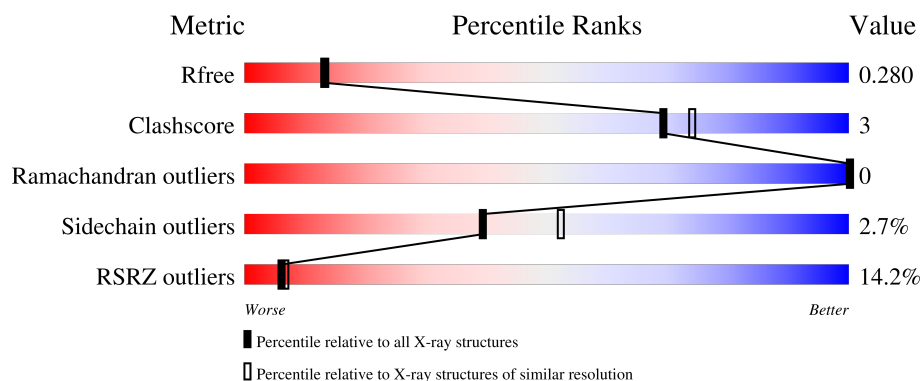
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.37 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	A	171	88% 9% ..
1	B	171	32% 84% 11% 5%
2	E	234	8% 81% 7% 11%
2	H	234	13% 81% 8% 11%
3	F	213	16% 78% 9% 10%

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Mol	Chain	Length	Quality of chain
3	L	213	10% 79% 12% • 6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8972 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 Lysosome-associated membrane glycoprotein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	168	Total	C	N	O	S	0	0	0
			1296	799	229	257	11			
1	B	162	Total	C	N	O	S	0	0	0
			1249	773	221	244	11			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25	GLY	-	expression tag	UNP P11279
A	26	SER	-	expression tag	UNP P11279
A	27	HIS	-	expression tag	UNP P11279
A	28	MET	-	expression tag	UNP P11279
B	25	GLY	-	expression tag	UNP P11279
B	26	SER	-	expression tag	UNP P11279
B	27	HIS	-	expression tag	UNP P11279
B	28	MET	-	expression tag	UNP P11279

- Molecule 2 is a protein called Fab B Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	208	Total	C	N	O	S	0	0	0
			1562	996	255	306	5			
2	H	208	Total	C	N	O	S	0	0	0
			1566	998	256	307	5			

- Molecule 3 is a protein called Fab B Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	192	Total	C	N	O	S	0	0	0
			1473	917	246	304	6			
3	L	201	Total	C	N	O	S	0	0	0
			1557	976	260	315	6			

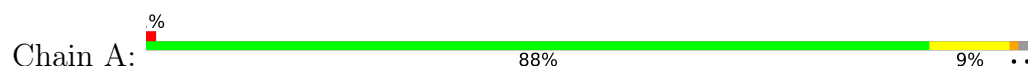
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	66	Total 66	O 66	0	0
4	B	18	Total 18	O 18	0	0
4	E	22	Total 22	O 22	0	0
4	F	32	Total 32	O 32	0	0
4	H	69	Total 69	O 69	0	0
4	L	62	Total 62	O 62	0	0

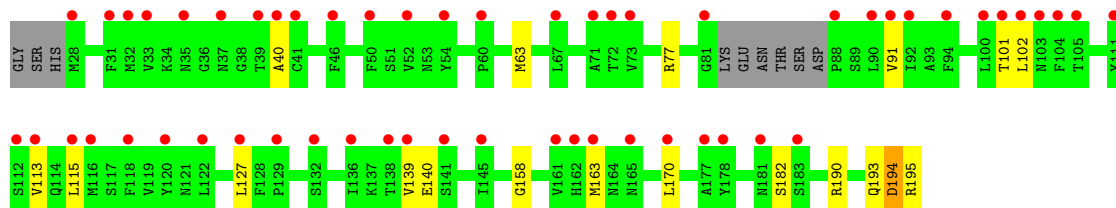
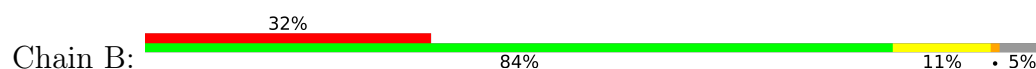
3 Residue-property plots [i](#)

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.

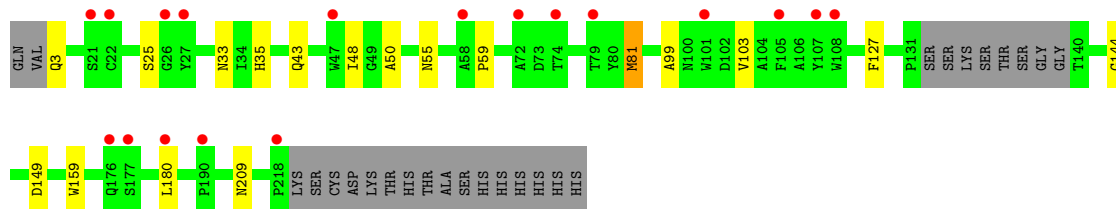
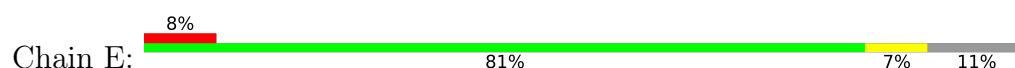
- Molecule 1: Lysosome-associated membrane glycoprotein 1



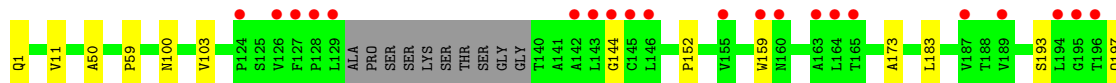
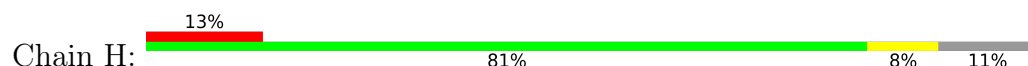
- Molecule 1: Lysosome-associated membrane glycoprotein 1

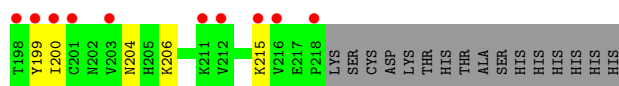


- Molecule 2: Fab B Heavy Chain

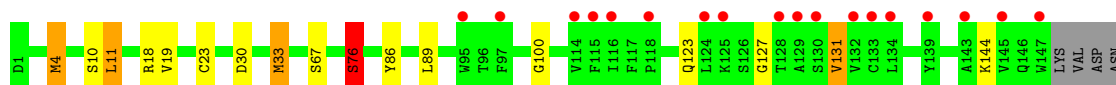
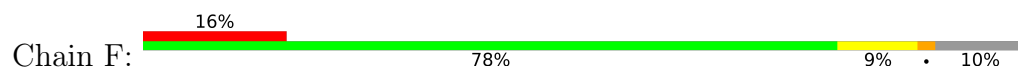


- Molecule 2: Fab B Heavy Chain

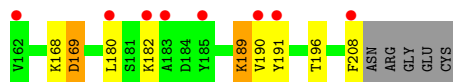
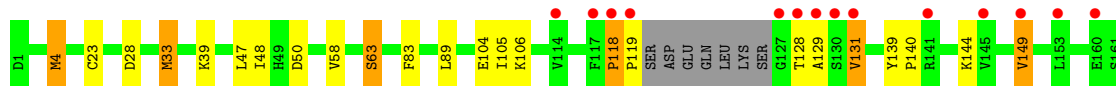
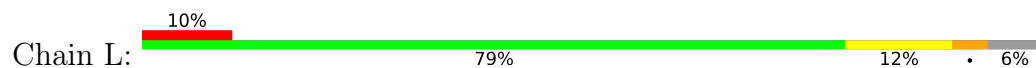




• Molecule 3: Fab B Light Chain



• Molecule 3: Fab B Light Chain



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	149.93Å 93.68Å 108.01Å 90.00° 115.93° 90.00°	Depositor
Resolution (Å)	72.08 – 2.37 72.08 – 2.37	Depositor EDS
% Data completeness (in resolution range)	99.0 (72.08-2.37) 98.9 (72.08-2.37)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.37Å)	Xtriage
Refinement program	BUSTER 2.11.8	Depositor
R, R_{free}	0.251 , 0.290 (Not available) , 0.280	Depositor DCC
R_{free} test set	2767 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	49.8	Xtriage
Anisotropy	0.388	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 34.7	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	0.93	EDS
Total number of atoms	8972	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% 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	A	0.78	0/1317	1.08	3/1780 (0.2%)
1	B	0.69	0/1269	1.04	2/1713 (0.1%)
2	E	0.68	1/1603 (0.1%)	0.98	0/2190
2	H	0.77	0/1606	1.00	1/2193 (0.0%)
3	F	0.72	2/1502 (0.1%)	0.98	2/2039 (0.1%)
3	L	0.82	3/1591 (0.2%)	1.07	7/2161 (0.3%)
All	All	0.75	6/8888 (0.1%)	1.02	15/12076 (0.1%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	33	MET	SD-CE	-8.31	1.58	1.79
3	F	33	MET	SD-CE	-7.92	1.59	1.79
3	L	128	THR	CA-C	5.40	1.59	1.52
3	L	4	MET	SD-CE	-5.18	1.66	1.79
3	F	4	MET	SD-CE	-5.13	1.66	1.79
2	E	81	MET	SD-CE	-5.06	1.67	1.79

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	144	ASP	CA-CB-CG	7.86	120.46	112.60
3	L	169	ASP	CA-CB-CG	7.41	120.01	112.60
3	L	118	PRO	O-C-N	6.98	124.43	121.15
1	B	194	ASP	CA-C-N	5.84	132.21	121.70
1	B	194	ASP	C-N-CA	5.84	132.21	121.70
3	L	169	ASP	CB-CA-C	-5.79	101.17	110.79
2	H	100	ASN	N-CA-C	-5.70	98.66	110.80
3	F	76	SER	N-CA-C	5.56	117.15	111.14
3	F	30	ASP	CA-CB-CG	5.45	118.05	112.60
1	A	194	ASP	CA-C-N	5.38	131.38	121.70
1	A	194	ASP	C-N-CA	5.38	131.38	121.70
3	L	28	ASP	CA-CB-CG	5.13	117.73	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	140	PRO	CA-C-N	5.13	127.40	120.38
3	L	140	PRO	C-N-CA	5.13	127.40	120.38
3	L	129	ALA	N-CA-C	5.10	117.80	109.59

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1296	0	1260	7	1
1	B	1249	0	1219	9	1
2	E	1562	0	1529	11	0
2	H	1566	0	1537	9	0
3	F	1473	0	1420	13	0
3	L	1557	0	1505	14	0
4	A	66	0	0	0	0
4	B	18	0	0	0	0
4	E	22	0	0	0	0
4	F	32	0	0	0	0
4	H	69	0	0	0	0
4	L	62	0	0	2	0
All	All	8972	0	8470	60	1

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 3.

All (60) 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:E:48:ILE:HD13	2:E:81:MET:HE1	1.68	0.75
3:F:192:ALA:HA	3:F:207:SER:HB3	1.74	0.69
3:F:11:LEU:HD13	3:F:19:VAL:HG13	1.78	0.66
3:L:119:PRO:HG3	3:L:131:VAL:HB	1.77	0.66
1:A:158:GLY:HA2	1:A:170:LEU:O	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:63:SER:HB3	4:L:318:HOH:O	2.01	0.59
3:F:18:ARG:HH21	3:F:76:SER:HG	1.50	0.58
1:B:158:GLY:HA2	1:B:170:LEU:O	2.03	0.58
3:F:127:GLY:HA2	3:F:182:LYS:HD2	1.86	0.58
1:B:63:MET:HG2	1:B:127:LEU:O	2.09	0.53
2:E:48:ILE:HD13	2:E:81:MET:CE	2.37	0.53
2:E:3:GLN:OE1	2:E:25:SER:OG	2.26	0.52
3:L:118:PRO:HB3	3:L:191:TYR:OH	2.10	0.52
2:E:103:VAL:HG11	3:F:89:LEU:HD23	1.92	0.52
3:F:192:ALA:HA	3:F:207:SER:CB	2.38	0.52
2:H:204:ASN:HD21	2:H:206:LYS:HE3	1.75	0.51
2:E:35:HIS:CD2	2:E:99:ALA:HB3	2.46	0.51
2:E:33:ASN:HB2	2:E:99:ALA:O	2.11	0.50
2:E:48:ILE:HG21	2:E:81:MET:HE3	1.93	0.50
1:B:115:LEU:HD11	1:B:140:GLU:HB3	1.95	0.48
1:A:63:MET:HG2	1:A:127:LEU:O	2.13	0.47
1:B:102:LEU:HD22	1:B:113:VAL:HG11	1.97	0.47
2:E:50:ALA:O	2:E:59:PRO:HD2	2.16	0.46
1:A:195:ARG:HA	1:A:195:ARG:HD3	1.70	0.46
1:A:40:ALA:HB3	1:A:182:SER:HA	1.98	0.46
1:B:77:ARG:HD3	1:B:91:VAL:HG11	1.97	0.46
3:F:180:LEU:HD13	3:F:184:ASP:HB2	1.98	0.46
1:B:40:ALA:HB3	1:B:182:SER:HA	1.97	0.46
2:H:50:ALA:O	2:H:59:PRO:HD2	2.16	0.45
2:H:144:GLY:HA2	2:H:159:TRP:CZ2	2.51	0.45
2:H:173:ALA:HA	2:H:183:LEU:HB3	1.98	0.45
2:E:149:ASP:HB3	2:E:180:LEU:HD13	1.99	0.45
3:F:4:MET:HE3	3:F:23:CYS:SG	2.57	0.45
1:B:139:VAL:HB	1:B:163:MET:HB3	1.99	0.44
3:L:144:LYS:HB3	3:L:196:THR:HB	1.99	0.44
3:L:149:VAL:HA	3:L:190:VAL:O	2.18	0.44
1:B:194:ASP:O	1:B:195:ARG:HB2	2.18	0.44
3:F:144:LYS:HB3	3:F:196:THR:HB	1.99	0.44
3:L:47:LEU:HA	3:L:58:VAL:HG21	2.01	0.43
2:E:144:GLY:HA2	2:E:159:TRP:CZ2	2.54	0.43
3:L:83:PHE:HB3	3:L:105:ILE:HG12	2.01	0.43
2:H:200:ILE:HG12	2:H:215:LYS:HG2	2.01	0.42
3:L:4:MET:HE3	3:L:23:CYS:SG	2.60	0.42
3:F:131:VAL:HG13	3:F:178:LEU:HB3	2.01	0.42
3:L:4:MET:CE	3:L:33:MET:HE1	2.49	0.42
3:L:106:LYS:HA	3:L:139:TYR:OH	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:4:MET:CE	3:F:33:MET:HE1	2.50	0.42
2:E:127:PHE:CE1	3:F:123:GLN:HA	2.54	0.42
2:H:11:VAL:HG21	2:H:152:PRO:HG3	2.01	0.42
2:H:199:TYR:O	2:H:215:LYS:HA	2.20	0.41
1:A:139:VAL:HB	1:A:163:MET:HB3	2.01	0.41
3:L:47:LEU:C	3:L:48:ILE:HG13	2.45	0.41
1:A:105:THR:HG22	1:A:114:GLN:HB2	2.03	0.41
2:H:193:SER:OG	2:H:197:GLN:HB3	2.21	0.41
3:L:39:LYS:HE2	4:L:302:HOH:O	2.20	0.41
3:L:189:LYS:HD2	3:L:190:VAL:HG23	2.02	0.40
1:B:91:VAL:HG13	1:B:101:THR:HG22	2.03	0.40
1:A:124:ASP:OD1	1:A:124:ASP:C	2.64	0.40
2:H:103:VAL:HG11	3:L:89:LEU:HD23	2.02	0.40
3:F:86:TYR:O	3:F:100:GLY:HA2	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:GLU:OE1	1:B:190:ARG:NH2[3_455]	2.14	0.06

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	166/171 (97%)	161 (97%)	5 (3%)	0	100	100
1	B	158/171 (92%)	155 (98%)	3 (2%)	0	100	100
2	E	204/234 (87%)	195 (96%)	9 (4%)	0	100	100
2	H	204/234 (87%)	197 (97%)	7 (3%)	0	100	100
3	F	186/213 (87%)	179 (96%)	7 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	L	197/213 (92%)	190 (96%)	7 (4%)	0	100	100
All	All	1115/1236 (90%)	1077 (97%)	38 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	A	148/150 (99%)	144 (97%)	4 (3%)	39	52
1	B	142/150 (95%)	141 (99%)	1 (1%)	76	86
2	E	175/198 (88%)	172 (98%)	3 (2%)	53	68
2	H	176/198 (89%)	175 (99%)	1 (1%)	78	87
3	F	169/188 (90%)	162 (96%)	7 (4%)	27	36
3	L	177/188 (94%)	166 (94%)	11 (6%)	16	19
All	All	987/1072 (92%)	960 (97%)	27 (3%)	39	52

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	124	ASP
1	A	180	SER
1	A	193	GLN
1	A	195	ARG
1	B	193	GLN
2	E	43	GLN
2	E	55	ASN
2	E	209	ASN
3	F	10	SER
3	F	11	LEU
3	F	67	SER
3	F	76	SER
3	F	131	VAL

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Mol	Chain	Res	Type
3	F	180	LEU
3	F	205	THR
2	H	1	GLN
3	L	50	ASP
3	L	63	SER
3	L	104	GLU
3	L	131	VAL
3	L	149	VAL
3	L	168	LYS
3	L	169	ASP
3	L	180	LEU
3	L	182	LYS
3	L	189	LYS
3	L	208	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	ASN
1	A	76	ASN
1	B	62	ASN
1	B	76	ASN
1	B	114	GLN
2	E	43	GLN
3	F	123	GLN
2	H	55	ASN
2	H	197	GLN
2	H	204	ASN
2	H	209	ASN

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 [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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	168/171 (98%)	0.17	1 (0%) 85 89	31, 46, 66, 83	0
1	B	162/171 (94%)	1.58	55 (33%) 1 0	42, 82, 102, 108	0
2	E	208/234 (88%)	0.92	18 (8%) 16 18	48, 70, 88, 99	0
2	H	208/234 (88%)	0.68	31 (14%) 5 6	30, 46, 93, 102	0
3	F	192/213 (90%)	1.03	35 (18%) 3 3	40, 68, 120, 126	0
3	L	201/213 (94%)	0.56	22 (10%) 10 12	33, 48, 85, 108	0
All	All	1139/1236 (92%)	0.81	162 (14%) 6 7	30, 60, 101, 126	0

All (162) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	199	TYR	6.2
2	H	216	VAL	6.0
2	H	128	PRO	5.8
3	L	127	GLY	5.1
1	B	161	VAL	4.8
1	B	104	PHE	4.7
3	F	114	VAL	4.6
2	H	126	VAL	4.6
2	H	159	TRP	4.5
2	H	145	CYS	4.5
3	L	191	TYR	4.4
2	H	189	VAL	4.4
1	B	102	LEU	4.2
2	H	127	PHE	4.2
2	H	129	LEU	4.0
1	B	90	LEU	4.0
1	B	81	GLY	3.9
3	F	195	VAL	3.9
2	H	144	GLY	3.8

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Mol	Chain	Res	Type	RSRZ
3	L	153	LEU	3.8
3	F	116	ILE	3.8
3	L	162	VAL	3.8
1	B	92	ILE	3.7
1	B	113	VAL	3.6
3	F	132	VAL	3.6
2	H	212	VAL	3.6
2	E	107	TYR	3.6
3	F	192	ALA	3.6
2	H	164	LEU	3.6
3	L	180	LEU	3.6
3	F	204	VAL	3.5
3	F	179	THR	3.5
3	F	130	SER	3.5
3	L	185	TYR	3.5
1	B	73	VAL	3.4
1	B	88	PRO	3.4
2	H	143	LEU	3.4
3	F	124	LEU	3.3
3	F	162	VAL	3.3
2	H	218	PRO	3.3
1	B	100	LEU	3.2
3	F	95	TRP	3.2
1	B	116	MET	3.2
2	H	198	THR	3.1
1	B	111	TYR	3.1
1	B	37	ASN	3.1
1	B	28	MET	3.1
3	L	208	PHE	3.0
3	F	205	THR	3.0
2	H	163	ALA	3.0
2	E	101	TRP	3.0
2	H	200	ILE	3.0
3	L	128	THR	3.0
2	E	218	PRO	2.9
1	B	52	VAL	2.9
1	B	71	ALA	2.9
1	B	50	PHE	2.8
2	H	155	VAL	2.8
1	B	41	CYS	2.8
2	H	142	ALA	2.8
1	B	139	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	177	ALA	2.7
1	B	31	PHE	2.7
1	B	118	PHE	2.7
2	H	195	GLY	2.7
3	L	130	SER	2.7
2	H	215	LYS	2.7
2	E	105	PHE	2.7
1	B	178	TYR	2.7
3	F	198	GLN	2.7
3	F	145	VAL	2.7
1	B	145	ILE	2.7
3	F	207	SER	2.7
2	H	146	LEU	2.7
2	E	190	PRO	2.6
3	F	134	LEU	2.6
3	L	183	ALA	2.6
1	B	72	THR	2.6
3	L	190	VAL	2.6
3	F	201	SER	2.6
1	B	115	LEU	2.5
3	L	129	ALA	2.5
1	B	46	PHE	2.5
1	B	60	PRO	2.5
1	B	138	THR	2.5
1	B	103	ASN	2.5
1	B	165	ASN	2.5
2	E	22	CYS	2.5
3	F	129	ALA	2.5
3	F	143	ALA	2.5
1	B	105	THR	2.5
2	H	165	THR	2.5
3	F	184	ASP	2.5
2	H	196	THR	2.5
1	B	33	VAL	2.5
1	B	91	VAL	2.5
2	E	177	SER	2.5
2	H	187	VAL	2.5
1	B	122	LEU	2.5
1	A	28	MET	2.4
1	B	162	HIS	2.4
3	F	118	PRO	2.4
1	B	94	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
2	E	108	TRP	2.4
3	F	180	LEU	2.4
1	B	101	THR	2.4
1	B	35	ASN	2.4
1	B	181	ASN	2.4
1	B	141	SER	2.4
3	L	145	VAL	2.4
3	L	149	VAL	2.4
3	L	160	GLU	2.4
2	E	26	GLY	2.3
1	B	112	SER	2.3
1	B	32	MET	2.3
1	B	127	LEU	2.3
3	F	139	TYR	2.3
3	F	147	TRP	2.3
1	B	163	MET	2.3
2	E	58	ALA	2.3
2	E	74	THR	2.3
2	E	79	THR	2.3
2	H	203	VAL	2.3
3	F	133	CYS	2.3
1	B	136	ILE	2.3
3	L	114	VAL	2.3
3	L	117	PHE	2.2
1	B	129	PRO	2.2
1	B	40	ALA	2.2
3	L	119	PRO	2.2
3	L	182	LYS	2.2
3	F	177	THR	2.2
1	B	54	TYR	2.2
2	E	47	TRP	2.2
2	H	194	LEU	2.2
3	F	115	PHE	2.2
3	L	118	PRO	2.2
2	H	211	LYS	2.2
3	F	128	THR	2.2
3	L	131	VAL	2.1
1	B	183	SER	2.1
2	E	21	SER	2.1
2	H	124	PRO	2.1
3	F	156	GLY	2.1
3	L	141	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	132	SER	2.1
3	F	196	THR	2.1
2	E	72	ALA	2.1
1	B	170	LEU	2.1
3	F	200	LEU	2.1
2	H	160	ASN	2.1
2	E	27	TYR	2.1
1	B	39	THR	2.0
2	H	201	CYS	2.0
3	F	193	CYS	2.0
1	B	67	LEU	2.0
1	B	120	TYR	2.0
3	F	97	PHE	2.0
3	F	161	SER	2.0
2	E	176	GLN	2.0
2	E	180	LEU	2.0
3	F	125	LYS	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.