



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 19, 2020 – 09:08 AM BST

PDB ID : 6LPC
Title : Crystal Structure of rat Munc18-1 with K332E/K333E mutation
Authors : Wang, X.P.; Gong, J.H.; Wang, S.; Zhu, L.; Yang, X.Y.; Xu, Y.Y.; Yang, X.F.;
Ma, C.
Deposited on : 2020-01-09
Resolution : 3.40 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	:	2.14.6
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)	:	2.14.6

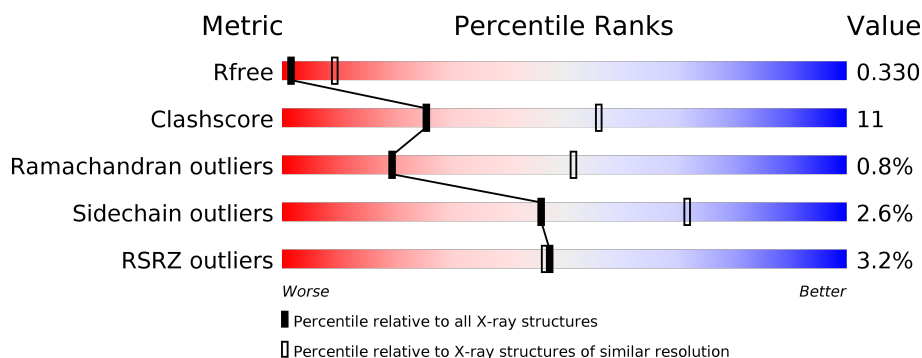
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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	1026 (3.48-3.32)
Clashscore	141614	1055 (3.48-3.32)
Ramachandran outliers	138981	1038 (3.48-3.32)
Sidechain outliers	138945	1038 (3.48-3.32)
RSRZ outliers	127900	2173 (3.50-3.30)

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	A	594	
1	B	594	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7114 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 Syntaxin-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	533	Total	C	N	O	S	0	0	0
			3814	2423	621	748	22			
1	B	467	Total	C	N	O	S	0	0	0
			3300	2081	548	654	17			

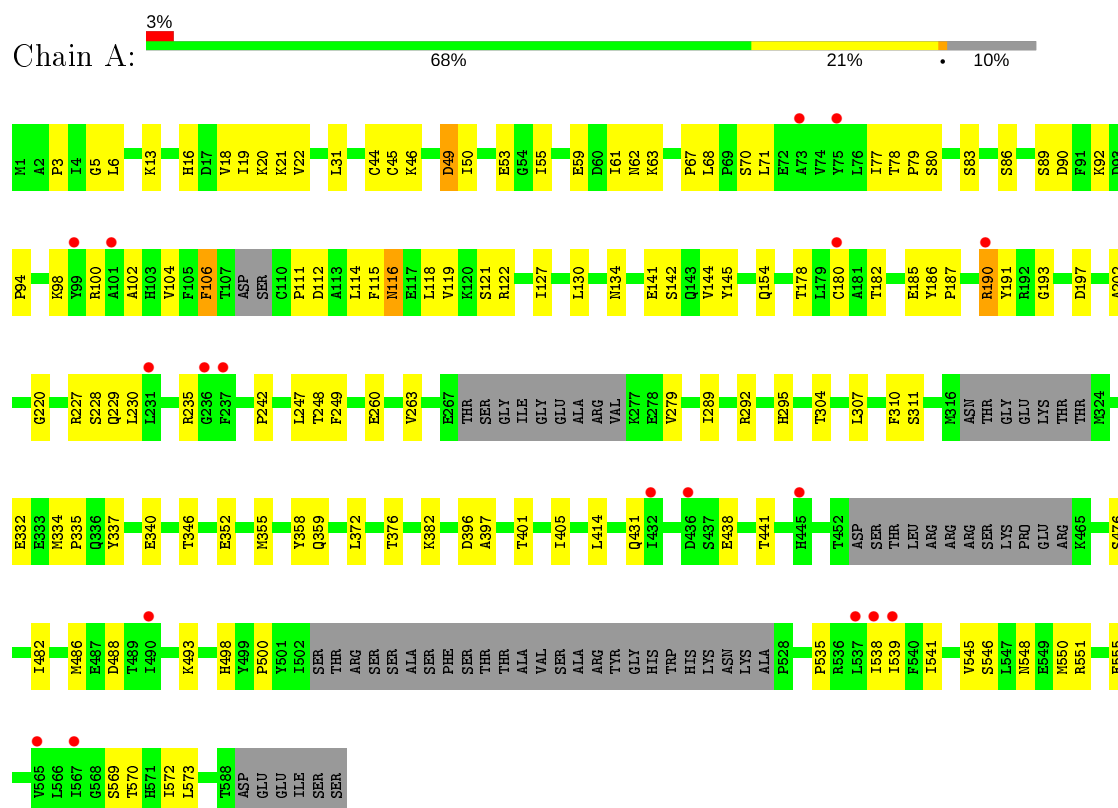
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	332	GLU	LYS	engineered mutation	UNP P61765
A	333	GLU	LYS	engineered mutation	UNP P61765
B	332	GLU	LYS	engineered mutation	UNP P61765
B	333	GLU	LYS	engineered mutation	UNP P61765

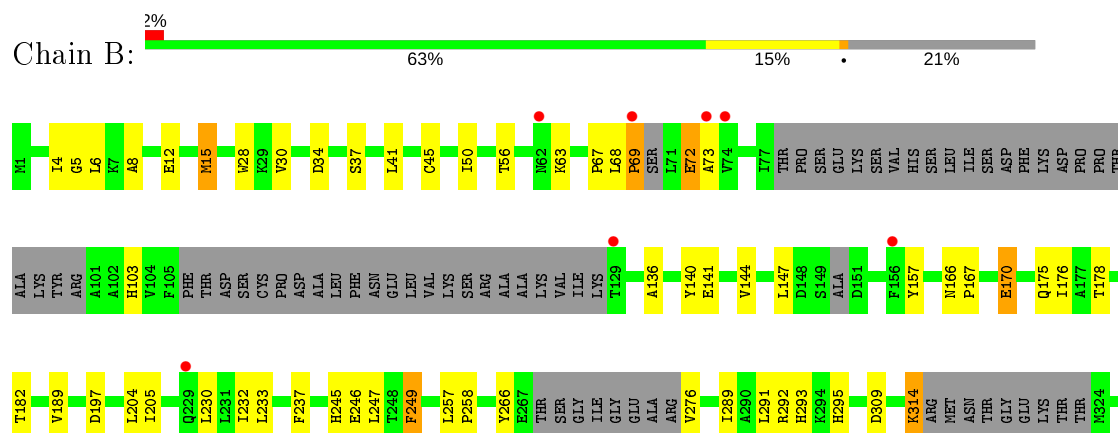
3 Residue-property plots [i](#)

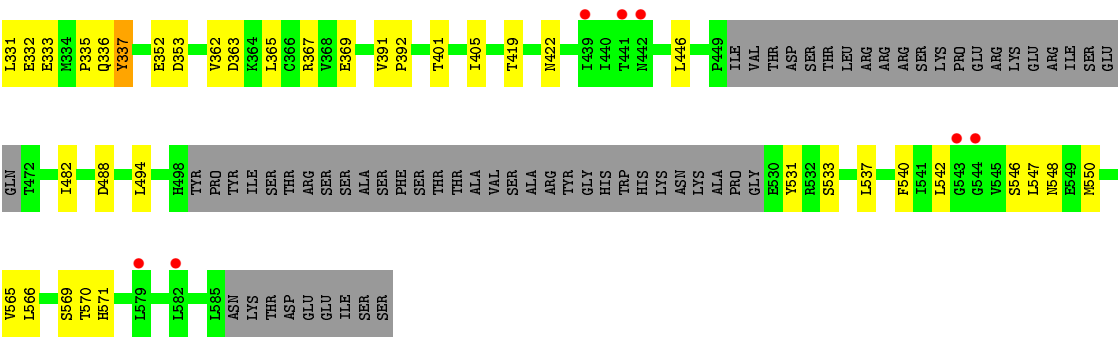
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Syntaxin-binding protein 1



• Molecule 1: Syntaxin-binding protein 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	80.47 Å 80.47 Å 247.66 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.07 – 3.40 49.07 – 3.40	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.07-3.40) 100.0 (49.07-3.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 3.40 Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.281 , 0.331 0.280 , 0.330	Depositor DCC
R_{free} test set	1104 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	126.2	Xtriage
Anisotropy	0.547	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 99.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.067 for h,-k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7114	wwPDB-VP
Average B, all atoms (Å ²)	135.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% 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	A	0.28	0/3887	0.49	0/5324
1	B	0.34	0/3354	0.53	1/4590 (0.0%)
All	All	0.31	0/7241	0.51	1/9914 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	69	PRO	N-CA-C	5.17	125.55	112.10

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	A	3814	0	3389	89	0
1	B	3300	0	2856	58	0
All	All	7114	0	6245	142	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 11.

All (142) 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:B:289:ILE:HA	1:B:292:ARG:HD2	1.61	0.82
1:B:6:LEU:HD21	1:B:566:LEU:HD21	1.61	0.81
1:A:68:LEU:HD13	1:A:70:SER:H	1.49	0.75
1:B:136:ALA:HB1	1:B:175:GLN:HG2	1.68	0.73
1:A:62:ASN:CB	1:A:86:SER:HB2	2.20	0.72
1:A:20:LYS:NZ	1:A:53:GLU:OE1	2.23	0.70
1:A:116:ASN:HA	1:A:119:VAL:HB	1.75	0.68
1:B:170:GLU:HG2	1:B:204:LEU:HD11	1.76	0.67
1:A:304:THR:HG21	1:B:314:LYS:HB2	1.76	0.67
1:A:13:LYS:HD2	1:A:182:THR:HG23	1.77	0.66
1:A:111:PRO:HB2	1:A:114:LEU:HB2	1.78	0.65
1:B:362:VAL:O	1:B:365:LEU:N	2.32	0.63
1:B:68:LEU:CB	1:B:69:PRO:HD2	2.29	0.63
1:A:551:ARG:NH1	1:A:555:GLU:OE2	2.32	0.63
1:B:4:ILE:HG12	1:B:5:GLY:H	1.65	0.62
1:A:5:GLY:HA3	1:A:142:SER:HA	1.81	0.61
1:B:257:LEU:HD21	1:B:353:ASP:HB3	1.82	0.61
1:B:245:HIS:ND1	1:B:369:GLU:OE2	2.30	0.61
1:A:279:VAL:HG21	1:A:346:THR:HG21	1.84	0.60
1:A:569:SER:OG	1:A:570:THR:N	2.35	0.60
1:A:334:MET:HG2	1:B:352:GLU:HB2	1.82	0.60
1:B:30:VAL:H	1:B:72:GLU:CB	2.14	0.60
1:B:419:THR:HG23	1:B:422:ASN:H	1.67	0.60
1:A:355:MET:O	1:A:359:GLN:N	2.32	0.59
1:B:141:GLU:HB2	1:B:144:VAL:HB	1.85	0.59
1:A:67:PRO:HB2	1:A:98:LYS:HA	1.85	0.58
1:A:67:PRO:CB	1:A:98:LYS:HA	2.33	0.58
1:A:68:LEU:HD12	1:A:71:LEU:HG	1.86	0.58
1:A:46:LYS:HB2	1:A:49:ASP:HB2	1.88	0.56
1:A:355:MET:HB2	1:B:331:LEU:HD12	1.87	0.55
1:B:249:PHE:HD2	1:B:291:LEU:HB3	1.72	0.55
1:A:134:ASN:O	1:A:134:ASN:ND2	2.40	0.55
1:A:396:ASP:OD1	1:A:397:ALA:N	2.40	0.54
1:A:248:THR:HG21	1:A:551:ARG:HD3	1.87	0.54
1:A:482:ILE:HD12	1:A:541:ILE:HD11	1.89	0.54
1:B:548:ASN:OD1	1:B:548:ASN:N	2.43	0.52
1:A:104:VAL:HG13	1:A:130:LEU:HD12	1.92	0.52
1:B:157:TYR:CE2	1:B:446:LEU:HA	2.45	0.52
1:B:332:GLU:O	1:B:335:PRO:HD2	2.10	0.51
1:A:45:CYS:CB	1:A:50:ILE:HD11	2.40	0.51
1:A:227:ARG:HB2	1:A:535:PRO:HB3	1.92	0.51
1:A:185:GLU:OE2	1:A:228:SER:OG	2.23	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:550:MET:HE2	1:A:572:ILE:HD12	1.93	0.50
1:A:551:ARG:HG2	1:A:555:GLU:HG3	1.91	0.50
1:B:531:TYR:CE2	1:B:533:SER:HB3	2.46	0.50
1:B:569:SER:OG	1:B:570:THR:N	2.37	0.50
1:A:68:LEU:CD1	1:A:70:SER:H	2.22	0.50
1:A:115:PHE:O	1:A:119:VAL:N	2.45	0.49
1:A:376:THR:HA	1:A:382:LYS:HA	1.93	0.48
1:A:178:THR:O	1:A:182:THR:OG1	2.23	0.48
1:A:45:CYS:HB3	1:A:50:ILE:HD11	1.96	0.48
1:A:46:LYS:HB2	1:A:49:ASP:CB	2.44	0.47
1:A:247:LEU:HD22	1:A:476:SER:HB3	1.96	0.47
1:B:488:ASP:O	1:B:494:LEU:N	2.48	0.47
1:A:77:ILE:HD11	1:A:106:PHE:HE1	1.79	0.47
1:B:230:LEU:HD21	1:B:232:ILE:HD11	1.97	0.47
1:B:546:SER:O	1:B:550:MET:HG3	2.14	0.47
1:B:73:ALA:O	1:B:103:HIS:HB2	2.15	0.47
1:A:114:LEU:O	1:A:118:LEU:N	2.48	0.47
1:A:438:GLU:O	1:A:441:THR:OG1	2.27	0.47
1:A:551:ARG:O	1:A:555:GLU:HG3	2.15	0.47
1:A:332:GLU:O	1:A:335:PRO:HD2	2.15	0.46
1:B:292:ARG:HG3	1:B:293:HIS:ND1	2.30	0.46
1:A:545:VAL:O	1:A:573:LEU:HB2	2.15	0.46
1:B:178:THR:O	1:B:182:THR:OG1	2.30	0.46
1:B:247:LEU:HD23	1:B:247:LEU:HA	1.77	0.46
1:A:141:GLU:HB2	1:A:144:VAL:HB	1.97	0.46
1:A:118:LEU:O	1:A:121:SER:OG	2.34	0.46
1:A:122:ARG:O	1:A:122:ARG:NE	2.48	0.46
1:A:235:ARG:HE	1:A:541:ILE:HG21	1.81	0.46
1:B:363:ASP:O	1:B:367:ARG:HG2	2.16	0.46
1:A:488:ASP:O	1:A:493:LYS:N	2.44	0.45
1:A:548:ASN:OD1	1:A:548:ASN:N	2.45	0.45
1:A:191:TYR:HB3	1:A:202:ALA:HB2	1.98	0.45
1:A:92:LYS:O	1:A:94:PRO:HD3	2.15	0.45
1:A:235:ARG:NH1	1:A:242:PRO:HG3	2.32	0.45
1:B:237:PHE:HZ	1:B:482:ILE:HB	1.80	0.45
1:B:189:VAL:HG23	1:B:230:LEU:HD22	1.98	0.45
1:B:8:ALA:O	1:B:12:GLU:N	2.44	0.45
1:A:352:GLU:OE2	1:B:335:PRO:HG3	2.16	0.45
1:A:5:GLY:HA3	1:A:142:SER:CA	2.46	0.45
1:B:333:GLU:O	1:B:337:TYR:N	2.45	0.45
1:A:401:THR:O	1:A:405:ILE:HG13	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:GLU:CB	1:A:63:LYS:HA	2.48	0.44
1:A:260:GLU:O	1:A:263:VAL:HG12	2.17	0.44
1:A:67:PRO:O	1:A:67:PRO:HG2	2.17	0.44
1:A:67:PRO:HB3	1:A:98:LYS:HA	2.00	0.44
1:A:115:PHE:CE1	1:A:130:LEU:HD21	2.53	0.44
1:A:78:THR:O	1:A:80:SER:N	2.51	0.44
1:A:16:HIS:O	1:A:21:LYS:HD3	2.18	0.43
1:A:186:TYR:HB2	1:A:220:GLY:HA3	1.99	0.43
1:B:15:MET:CE	1:B:50:ILE:HG22	2.48	0.43
1:B:247:LEU:HB3	1:B:295:HIS:NE2	2.33	0.43
1:A:546:SER:O	1:A:550:MET:HG3	2.19	0.43
1:B:15:MET:HE1	1:B:50:ILE:HG22	2.01	0.43
1:B:15:MET:HE1	1:B:50:ILE:HA	2.01	0.43
1:B:140:TYR:CD2	1:B:141:GLU:HG3	2.54	0.43
1:B:246:GLU:HG3	1:B:548:ASN:HB3	2.00	0.43
1:B:68:LEU:CB	1:B:69:PRO:CD	2.96	0.43
1:A:307:LEU:O	1:A:311:SER:OG	2.35	0.43
1:B:233:LEU:O	1:B:542:LEU:N	2.46	0.43
1:B:537:LEU:HD23	1:B:565:VAL:HG22	2.00	0.43
1:A:193:GLY:N	1:A:498:HIS:O	2.42	0.43
1:A:355:MET:CB	1:B:331:LEU:HD12	2.49	0.43
1:A:78:THR:C	1:A:80:SER:H	2.22	0.43
1:A:19:ILE:HD12	1:A:19:ILE:H	1.83	0.43
1:B:391:VAL:N	1:B:392:PRO:HD2	2.34	0.43
1:A:230:LEU:HD13	1:A:538:ILE:HB	2.00	0.43
1:A:372:LEU:O	1:A:414:LEU:HD13	2.18	0.43
1:B:28:TRP:CZ3	1:B:56:THR:HB	2.54	0.43
1:A:187:PRO:HB2	1:A:229:GLN:C	2.40	0.42
1:B:266:TYR:O	1:B:276:VAL:N	2.52	0.42
1:A:180:CYS:HB3	1:A:185:GLU:O	2.20	0.42
1:A:289:ILE:HA	1:A:292:ARG:HD3	2.02	0.42
1:A:31:LEU:HB2	1:A:55:ILE:HD13	2.00	0.42
1:A:112:ASP:N	1:A:112:ASP:OD2	2.53	0.42
1:A:115:PHE:HE1	1:A:130:LEU:HD21	1.85	0.42
1:A:310:PHE:CZ	1:A:340:GLU:HG2	2.55	0.42
1:A:541:ILE:O	1:A:569:SER:OG	2.29	0.42
1:B:34:ASP:H	1:B:37:SER:HG	1.64	0.42
1:A:6:LEU:HD12	1:A:145:TYR:HB3	2.02	0.42
1:B:166:ASN:N	1:B:167:PRO:HD2	2.35	0.42
1:A:86:SER:O	1:A:89:SER:N	2.52	0.42
1:B:247:LEU:HB3	1:B:295:HIS:CD2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:482:ILE:HD11	1:A:539:ILE:HG21	2.02	0.42
1:B:147:LEU:HD11	1:B:540:PHE:CE1	2.55	0.42
1:B:309:ASP:OD1	1:B:309:ASP:N	2.52	0.42
1:A:102:ALA:HB3	1:A:127:ILE:HA	2.02	0.41
1:B:362:VAL:O	1:B:363:ASP:C	2.57	0.41
1:A:18:VAL:O	1:A:22:VAL:HG22	2.20	0.41
1:A:83:SER:O	1:A:86:SER:OG	2.30	0.41
1:B:4:ILE:HA	1:B:8:ALA:HB2	2.03	0.41
1:A:3:PRO:HG2	1:A:44:CYS:O	2.20	0.41
1:B:401:THR:O	1:B:405:ILE:HG13	2.21	0.41
1:B:547:LEU:HG	1:B:550:MET:HE3	2.02	0.41
1:A:355:MET:HA	1:A:358:TYR:HB3	2.03	0.40
1:A:482:ILE:O	1:A:486:MET:HG3	2.21	0.40
1:B:176:ILE:HD12	1:B:205:ILE:HD11	2.03	0.40
1:A:190:ARG:HG2	1:A:500:PRO:O	2.21	0.40
1:A:247:LEU:HD23	1:A:295:HIS:CE1	2.56	0.40
1:B:41:LEU:HD12	1:B:45:CYS:HB2	2.03	0.40
1:A:191:TYR:CE2	1:A:500:PRO:HG2	2.57	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	A	521/594 (88%)	461 (88%)	56 (11%)	4 (1%)	19	51
1	B	449/594 (76%)	401 (89%)	44 (10%)	4 (1%)	17	49
All	All	970/1188 (82%)	862 (89%)	100 (10%)	8 (1%)	19	51

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	61	ILE
1	B	63	LYS
1	A	431	GLN
1	A	79	PRO
1	A	100	ARG
1	B	67	PRO
1	B	72	GLU
1	B	258	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	A	358/533 (67%)	349 (98%)	9 (2%)	47	72
1	B	298/533 (56%)	290 (97%)	8 (3%)	44	70
All	All	656/1066 (62%)	639 (97%)	17 (3%)	46	72

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	ASP
1	A	90	ASP
1	A	106	PHE
1	A	116	ASN
1	A	154	GLN
1	A	190	ARG
1	A	197	ASP
1	A	249	PHE
1	A	337	TYR
1	B	15	MET
1	B	170	GLU
1	B	197	ASP
1	B	249	PHE
1	B	314	LYS
1	B	336	GLN
1	B	337	TYR

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Mol	Chain	Res	Type
1	B	571	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	442	ASN
1	A	445	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 monosaccharides 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	533/594 (89%)	0.02	18 (3%) 45 44	30, 130, 157, 179	0
1	B	467/594 (78%)	0.04	14 (2%) 50 49	89, 138, 193, 244	0
All	All	1000/1188 (84%)	0.03	32 (3%) 47 46	30, 134, 182, 244	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	101	ALA	4.3
1	A	99	TYR	4.3
1	B	579	LEU	3.9
1	A	539	ILE	3.7
1	A	75	TYR	3.3
1	B	229	GLN	3.1
1	A	231	LEU	3.1
1	B	129	THR	3.1
1	A	538	ILE	3.0
1	B	543	GLY	2.9
1	B	156	PHE	2.8
1	A	565	VAL	2.6
1	B	73	ALA	2.5
1	A	567	ILE	2.5
1	A	490	ILE	2.5
1	A	236	GLY	2.4
1	A	537	LEU	2.4
1	A	180	CYS	2.4
1	A	237	PHE	2.4
1	A	436	ASP	2.3
1	B	441	THR	2.3
1	B	544	GLY	2.2
1	A	73	ALA	2.2
1	A	445	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	69	PRO	2.2
1	A	432	ILE	2.2
1	A	190	ARG	2.1
1	B	582	LEU	2.1
1	B	62	ASN	2.1
1	B	442	ASN	2.1
1	B	439	ILE	2.0
1	B	74	VAL	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 monosaccharides 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.