



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:31 PM UTC

PDB ID : 8FTU / pdb_00008ftu
Title : Crystal structure of the SNARE Use1 bound to Dsl1 complex subunits Sec39 and Dsl1, Revised Use1 structure
Authors : Travis, S.M.; Jeffrey, P.D.; Hughson, F.M.
Deposited on : 2023-01-13
Resolution : 5.73 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

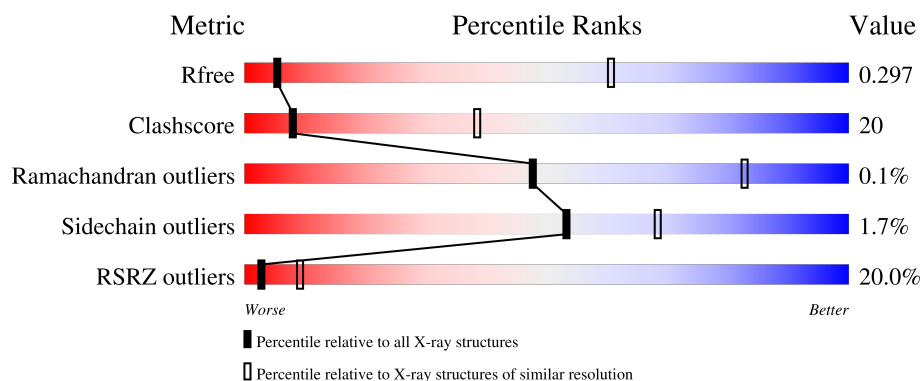
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 5.73 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	1110 (7.46-4.00)
Clashscore	190562	1171 (7.46-4.00)
Ramachandran outliers	187476	1005 (7.40-4.00)
Sidechain outliers	187428	1004 (7.50-3.98)
RSRZ outliers	180081	1103 (7.46-4.00)

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	A	699	23% 83% 7% 10%
2	B	113	15% 73% 8% 19%
3	C	306	8% 39% 50% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8254 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 Protein transport protein SEC39.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	627	Total	C	N	O	S	0	0	0
			5059	3255	817	969	18			

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-26	MET	-	expression tag	UNP Q6CWC7
A	-25	LYS	-	expression tag	UNP Q6CWC7
A	-24	HIS	-	expression tag	UNP Q6CWC7
A	-23	HIS	-	expression tag	UNP Q6CWC7
A	-22	HIS	-	expression tag	UNP Q6CWC7
A	-21	HIS	-	expression tag	UNP Q6CWC7
A	-20	HIS	-	expression tag	UNP Q6CWC7
A	-19	HIS	-	expression tag	UNP Q6CWC7
A	-18	HIS	-	expression tag	UNP Q6CWC7
A	-17	GLY	-	expression tag	UNP Q6CWC7
A	-16	ALA	-	expression tag	UNP Q6CWC7
A	-15	ALA	-	expression tag	UNP Q6CWC7
A	-14	GLY	-	expression tag	UNP Q6CWC7
A	-13	THR	-	expression tag	UNP Q6CWC7
A	-12	SER	-	expression tag	UNP Q6CWC7
A	-11	LEU	-	expression tag	UNP Q6CWC7
A	-10	TYR	-	expression tag	UNP Q6CWC7
A	-9	LYS	-	expression tag	UNP Q6CWC7
A	-8	LYS	-	expression tag	UNP Q6CWC7
A	-7	ALA	-	expression tag	UNP Q6CWC7
A	-6	GLY	-	expression tag	UNP Q6CWC7
A	-5	LEU	-	expression tag	UNP Q6CWC7
A	-4	VAL	-	expression tag	UNP Q6CWC7
A	-3	PRO	-	expression tag	UNP Q6CWC7
A	-2	ARG	-	expression tag	UNP Q6CWC7
A	-1	GLY	-	expression tag	UNP Q6CWC7
A	0	SER	-	expression tag	UNP Q6CWC7

- Molecule 2 is a protein called Vesicle transport protein USE1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	91	Total	C	N	O	0	0	0
			743	465	127	151			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	expression tag	UNP Q6CL08
B	2	GLY	-	expression tag	UNP Q6CL08
B	3	SER	-	expression tag	UNP Q6CL08

- Molecule 3 is a protein called Protein transport protein DSL1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	294	Total	C	N	O	S	0	0	0
			2452	1588	402	454	8			

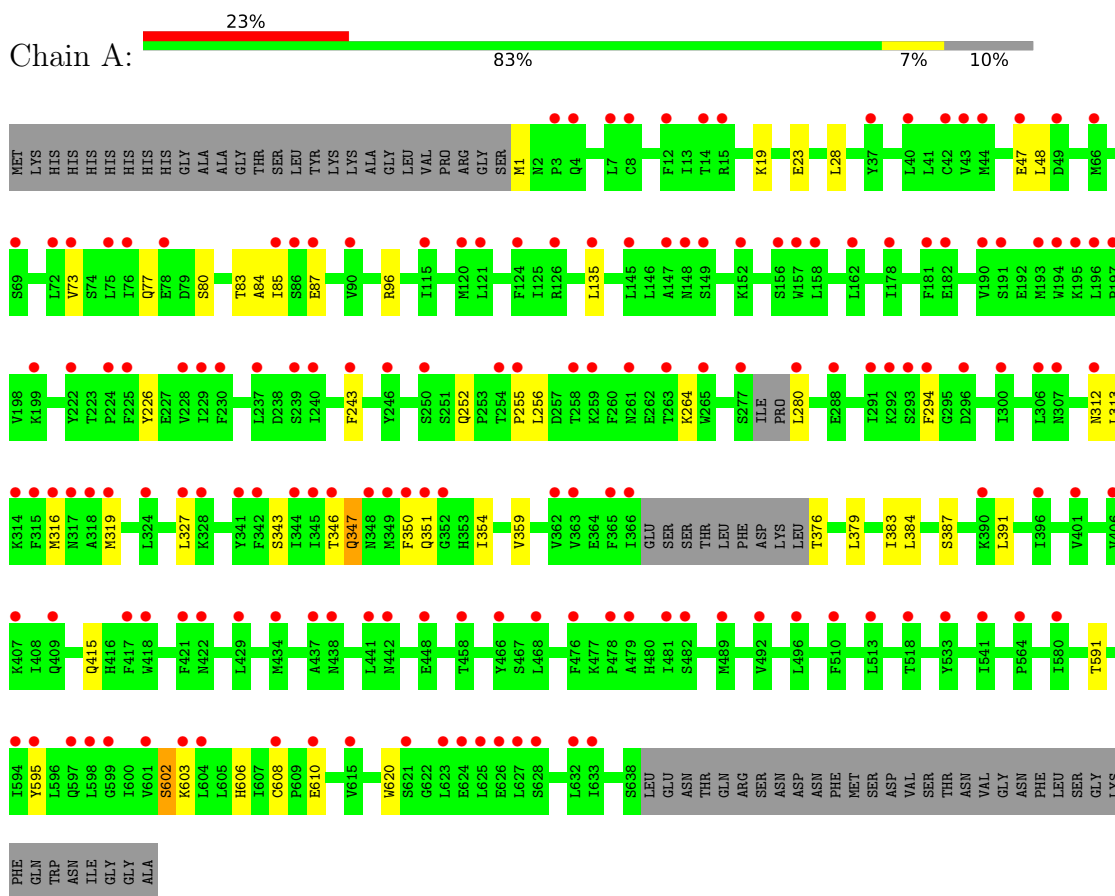
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	419	MET	-	expression tag	UNP Q6CUS2
C	420	GLY	-	expression tag	UNP Q6CUS2
C	421	SER	-	expression tag	UNP Q6CUS2
C	509	GLY	-	linker	UNP Q6CUS2
C	510	ASP	-	linker	UNP Q6CUS2
C	511	GLY	-	linker	UNP Q6CUS2
C	512	ASP	-	linker	UNP Q6CUS2
C	513	GLY	-	linker	UNP Q6CUS2

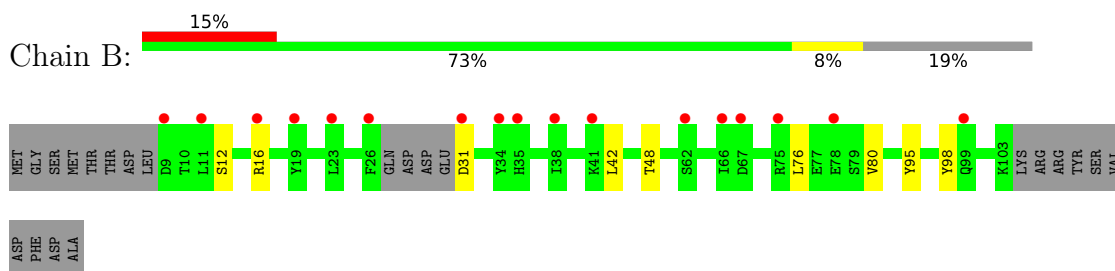
3 Residue-property plots [i](#)

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.

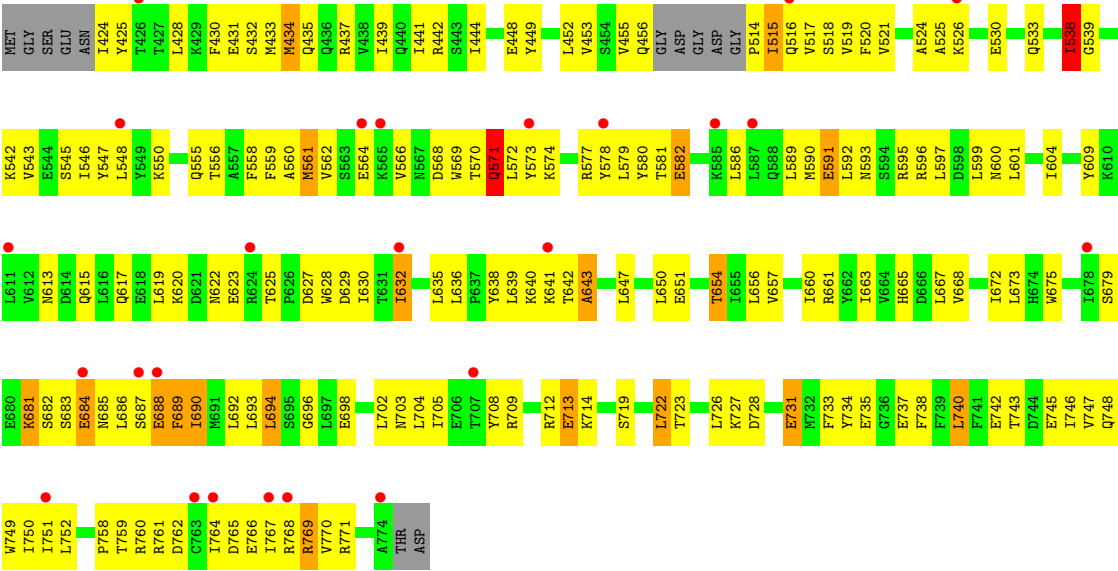
- Molecule 1: Protein transport protein SEC39



- Molecule 2: Vesicle transport protein USE1



- Molecule 3: Protein transport protein DSL1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	240.79Å 87.85Å 161.43Å 90.00° 107.61° 90.00°	Depositor
Resolution (Å)	29.60 – 5.73 29.60 – 5.73	Depositor EDS
% Data completeness (in resolution range)	74.5 (29.60-5.73) 73.9 (29.60-5.73)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.33 (at 5.67Å)	Xtriage
Refinement program	PHENIX 1.17.1-3660-000	Depositor
R, R_{free}	0.276 , 0.298 0.274 , 0.297	Depositor DCC
R_{free} test set	706 reflections (7.41%)	wwPDB-VP
Wilson B-factor (Å ²)	317.3	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 255.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	8254	wwPDB-VP
Average B, all atoms (Å ²)	271.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.59% 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	0.69	0/5159	1.19	4/6989 (0.1%)
2	B	0.69	0/754	1.21	1/1014 (0.1%)
3	C	0.83	0/2497	1.44	26/3378 (0.8%)
All	All	0.73	0/8410	1.27	31/11381 (0.3%)

There are no bond length outliers.

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	434	MET	N-CA-C	-8.38	101.19	111.33
3	C	689	PHE	N-CA-C	8.36	120.09	110.97
3	C	651	GLU	N-CA-C	7.40	121.60	111.39
1	A	77	GLN	OE1-CD-NE2	-7.12	115.48	122.60
3	C	681	LYS	CA-CB-CG	-6.33	101.44	114.10
2	B	31	ASP	CA-CB-CG	-6.32	106.28	112.60
3	C	654	THR	CA-C-N	6.28	128.60	120.56
3	C	654	THR	C-N-CA	6.28	128.60	120.56
3	C	684	GLU	CA-CB-CG	-6.23	101.63	114.10
3	C	538	ILE	CA-C-N	6.11	132.16	120.07
3	C	538	ILE	C-N-CA	6.11	132.16	120.07
3	C	561	MET	N-CA-C	5.76	117.56	111.28
1	A	347	GLN	OE1-CD-NE2	-5.71	116.89	122.60
3	C	769	ARG	CG-CD-NE	-5.67	99.53	112.00
3	C	444	ILE	N-CA-C	-5.57	96.84	108.88
3	C	547	TYR	CA-CB-CG	5.52	123.83	113.90
3	C	731	GLU	N-CA-CB	-5.41	101.94	110.06
3	C	632	ILE	N-CA-CB	-5.38	102.50	110.58
3	C	643	ALA	CA-C-N	5.38	126.46	119.78
3	C	643	ALA	C-N-CA	5.38	126.46	119.78
1	A	415	GLN	OE1-CD-NE2	-5.26	117.34	122.60
3	C	722	LEU	N-CA-C	5.25	117.75	111.71
3	C	571	GLN	CA-CB-CG	5.21	124.52	114.10
3	C	688	GLU	CA-C-N	5.20	127.52	120.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	688	GLU	C-N-CA	5.20	127.52	120.65
1	A	252	GLN	OE1-CD-NE2	-5.19	117.41	122.60
3	C	582	GLU	CA-CB-CG	5.16	124.42	114.10
3	C	737	GLU	CA-CB-CG	-5.15	103.79	114.10
3	C	690	ILE	O-C-N	5.12	127.06	121.94
3	C	433	MET	N-CA-C	-5.03	105.89	112.23
3	C	545	SER	N-CA-C	5.01	117.64	111.82

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	5059	0	5084	123	0
2	B	743	0	719	22	0
3	C	2452	0	2506	225	0
All	All	8254	0	8309	325	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 20.

All (325) 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:A:343:SER:HB3	1:A:383:ILE:CD1	1.15	1.60
1:A:343:SER:CB	1:A:383:ILE:HD13	1.22	1.58
1:A:602:SER:CB	3:C:560:ALA:HB1	1.51	1.40
1:A:343:SER:CB	1:A:383:ILE:HG21	1.50	1.38
1:A:343:SER:CA	1:A:383:ILE:HG21	1.50	1.38
1:A:347:GLN:HG2	1:A:387:SER:CB	1.58	1.33
1:A:343:SER:HB2	1:A:383:ILE:CG2	1.61	1.30
1:A:347:GLN:CG	1:A:387:SER:HB3	1.61	1.29
3:C:640:LYS:HE2	3:C:696:GLY:HA2	1.23	1.16
1:A:226:TYR:OH	1:A:255:PRO:CG	2.02	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:602:SER:CB	3:C:560:ALA:CB	2.30	1.08
1:A:603:LYS:HG3	3:C:564:GLU:HG3	1.23	1.08
3:C:425:TYR:HA	3:C:428:LEU:HD12	1.26	1.07
3:C:424:ILE:O	3:C:428:LEU:HG	1.53	1.06
1:A:602:SER:HB3	3:C:560:ALA:CB	1.86	1.06
1:A:350:PHE:HE1	1:A:359:VAL:HG22	1.17	1.04
1:A:603:LYS:CG	3:C:564:GLU:HG3	1.88	1.04
1:A:83:THR:OG1	2:B:95:TYR:CZ	2.15	0.99
1:A:343:SER:CA	1:A:383:ILE:CG2	2.42	0.97
1:A:343:SER:CB	1:A:383:ILE:CG2	2.28	0.97
1:A:343:SER:HB2	1:A:383:ILE:HG21	1.17	0.97
3:C:690:ILE:O	3:C:693:LEU:HB3	1.65	0.95
1:A:347:GLN:HE21	1:A:383:ILE:HG23	1.29	0.95
1:A:620:TRP:CZ3	3:C:556:THR:CG2	2.49	0.94
1:A:343:SER:HA	1:A:383:ILE:HG21	1.49	0.94
1:A:602:SER:HB2	3:C:560:ALA:HB1	1.46	0.94
3:C:758:PRO:HA	3:C:761:ARG:HB2	1.52	0.92
1:A:84:ALA:CB	2:B:95:TYR:HB2	2.01	0.91
1:A:350:PHE:CE1	1:A:359:VAL:HG22	2.06	0.91
1:A:346:THR:O	1:A:350:PHE:HD2	1.54	0.90
1:A:603:LYS:HE2	3:C:564:GLU:HB3	1.53	0.90
3:C:613:ASN:O	3:C:617:GLN:N	2.06	0.89
1:A:343:SER:HB3	1:A:383:ILE:CG1	2.01	0.89
3:C:521:VAL:HG11	3:C:578:TYR:CD2	2.08	0.89
3:C:638:TYR:O	3:C:643:ALA:N	2.06	0.89
2:B:16:ARG:CZ	2:B:76:LEU:HB2	2.03	0.88
1:A:226:TYR:OH	1:A:255:PRO:HG2	1.72	0.88
1:A:620:TRP:CZ3	3:C:556:THR:HG23	2.10	0.87
1:A:243:PHE:CE1	1:A:280:LEU:HD11	2.09	0.87
3:C:580:TYR:CE1	3:C:590:MET:HE3	2.11	0.86
1:A:343:SER:C	1:A:383:ILE:HG21	2.00	0.86
1:A:602:SER:OG	3:C:560:ALA:HB1	1.74	0.86
3:C:430:PHE:O	3:C:434:MET:HG2	1.74	0.86
3:C:639:LEU:HA	3:C:643:ALA:HB3	1.59	0.85
1:A:343:SER:CB	1:A:383:ILE:CG1	2.54	0.85
3:C:719:SER:HA	3:C:722:LEU:HD12	1.60	0.83
1:A:343:SER:CA	1:A:383:ILE:HD13	2.07	0.83
2:B:16:ARG:NH1	2:B:76:LEU:HB2	1.94	0.82
3:C:448:GLU:HB3	3:C:520:PHE:HB3	1.62	0.81
1:A:84:ALA:HB2	2:B:95:TYR:HB2	1.62	0.81
1:A:602:SER:HB3	3:C:560:ALA:HB3	1.63	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:580:TYR:CD2	3:C:590:MET:HG2	2.16	0.81
3:C:751:ILE:HA	3:C:760:ARG:NH1	1.96	0.81
3:C:555:GLN:HB3	3:C:589:LEU:HG	1.60	0.81
1:A:264:LYS:HE2	1:A:294:PHE:CZ	2.16	0.80
3:C:546:ILE:HD12	3:C:550:LYS:HE3	1.64	0.80
3:C:425:TYR:O	3:C:428:LEU:HB2	1.81	0.79
1:A:347:GLN:HE21	1:A:383:ILE:CG2	1.97	0.78
1:A:343:SER:HA	1:A:383:ILE:CG2	2.12	0.78
1:A:83:THR:HG21	2:B:95:TYR:CE1	2.19	0.77
3:C:569:TRP:CZ2	3:C:600:ASN:HB3	2.19	0.77
3:C:425:TYR:HA	3:C:428:LEU:CD1	2.11	0.77
3:C:580:TYR:CG	3:C:590:MET:HG2	2.21	0.76
1:A:226:TYR:OH	1:A:255:PRO:HG3	1.83	0.76
1:A:343:SER:HB2	1:A:383:ILE:HG23	1.63	0.76
3:C:647:LEU:HB3	3:C:656:LEU:HD21	1.68	0.75
3:C:726:LEU:HD23	3:C:759:THR:HG22	1.68	0.75
3:C:617:GLN:HE22	3:C:620:LYS:HD2	1.51	0.75
1:A:350:PHE:HD1	1:A:354:ILE:HD13	1.51	0.74
3:C:580:TYR:CD1	3:C:590:MET:HG2	2.21	0.74
1:A:350:PHE:CZ	1:A:384:LEU:HD11	2.23	0.74
3:C:615:GLN:HA	3:C:630:ILE:HD12	1.70	0.74
3:C:620:LYS:HE2	3:C:675:TRP:CE2	2.23	0.74
3:C:521:VAL:HG11	3:C:578:TYR:CE2	2.23	0.73
3:C:619:LEU:HD12	3:C:686:LEU:HD23	1.67	0.73
1:A:87:GLU:O	2:B:98:TYR:CE1	2.41	0.73
1:A:84:ALA:HB2	2:B:95:TYR:CB	2.19	0.73
3:C:733:PHE:HB2	3:C:738:PHE:CE2	2.24	0.73
3:C:568:ASP:HB3	3:C:571:GLN:HB2	1.69	0.72
3:C:566:VAL:HB	3:C:572:LEU:HD13	1.71	0.72
1:A:591:THR:HG22	1:A:595:TYR:HE2	1.54	0.71
3:C:690:ILE:O	3:C:693:LEU:CB	2.38	0.71
1:A:350:PHE:CZ	1:A:384:LEU:CD1	2.74	0.71
3:C:586:LEU:HB3	3:C:589:LEU:HD12	1.72	0.70
1:A:83:THR:CG2	2:B:95:TYR:CE1	2.74	0.70
1:A:47:GLU:OE1	1:A:135:LEU:HB2	1.92	0.70
3:C:521:VAL:CG1	3:C:578:TYR:CE2	2.74	0.70
3:C:569:TRP:CZ2	3:C:600:ASN:CB	2.74	0.70
3:C:628:TRP:O	3:C:632:ILE:HG13	1.92	0.69
3:C:668:VAL:O	3:C:672:ILE:HG13	1.93	0.69
3:C:580:TYR:CE2	3:C:590:MET:HG2	2.27	0.68
3:C:746:ILE:HD11	3:C:771:ARG:NH1	2.08	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:GLU:OE1	1:A:135:LEU:HD22	1.94	0.68
3:C:628:TRP:HH2	3:C:685:ASN:O	1.76	0.68
1:A:620:TRP:HZ3	3:C:556:THR:CG2	2.05	0.68
1:A:350:PHE:HD1	1:A:354:ILE:CD1	2.07	0.68
1:A:343:SER:CB	1:A:383:ILE:CD1	2.07	0.67
1:A:350:PHE:CD1	1:A:354:ILE:HD13	2.29	0.67
3:C:521:VAL:HG12	3:C:578:TYR:HE2	1.59	0.67
1:A:350:PHE:CE2	1:A:384:LEU:HD12	2.29	0.67
1:A:347:GLN:NE2	1:A:383:ILE:HG23	2.06	0.67
3:C:580:TYR:CG	3:C:590:MET:CG	2.78	0.66
3:C:668:VAL:HG21	3:C:693:LEU:HD21	1.77	0.66
3:C:638:TYR:CE2	3:C:643:ALA:HB2	2.30	0.65
3:C:591:GLU:HB3	3:C:595:ARG:HH12	1.60	0.65
3:C:708:TYR:O	3:C:712:ARG:HG3	1.97	0.65
3:C:762:ASP:O	3:C:766:GLU:HB2	1.97	0.65
3:C:521:VAL:HG11	3:C:578:TYR:HD2	1.60	0.65
3:C:733:PHE:CZ	3:C:771:ARG:HG2	2.32	0.65
1:A:602:SER:O	3:C:561:MET:HG2	1.98	0.64
3:C:731:GLU:O	3:C:735:GLU:HG3	1.98	0.64
3:C:733:PHE:HZ	3:C:771:ARG:HG2	1.62	0.64
1:A:47:GLU:OE1	1:A:135:LEU:CB	2.46	0.64
1:A:346:THR:O	1:A:350:PHE:CD2	2.44	0.64
3:C:580:TYR:CE1	3:C:590:MET:HG2	2.33	0.64
3:C:657:VAL:HG21	3:C:705:ILE:HG21	1.78	0.64
3:C:530:GLU:HA	3:C:533:GLN:HB2	1.79	0.63
1:A:620:TRP:HZ3	3:C:556:THR:HG23	1.60	0.63
3:C:432:SER:O	3:C:435:GLN:HB2	1.99	0.63
1:A:83:THR:OG1	2:B:95:TYR:CE2	2.44	0.63
1:A:312:ASN:HB3	1:A:327:LEU:HD21	1.81	0.63
3:C:684:GLU:HA	3:C:687:SER:HB3	1.81	0.62
3:C:734:TYR:HE1	3:C:770:VAL:HG21	1.64	0.62
1:A:48:LEU:O	2:B:48:THR:HG21	1.99	0.62
1:A:343:SER:HB2	1:A:383:ILE:CG1	2.25	0.62
3:C:596:ARG:O	3:C:600:ASN:ND2	2.29	0.62
3:C:521:VAL:O	3:C:524:ALA:HB3	1.99	0.62
3:C:515:ILE:HG13	3:C:516:GLN:N	2.13	0.62
1:A:243:PHE:CE1	1:A:280:LEU:CD1	2.83	0.62
1:A:603:LYS:CE	3:C:564:GLU:HB3	2.29	0.62
1:A:610:GLU:HB3	3:C:435:GLN:HG2	1.83	0.61
3:C:690:ILE:O	3:C:693:LEU:N	2.33	0.61
3:C:580:TYR:CZ	3:C:590:MET:HG2	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:766:GLU:HA	3:C:769:ARG:NH1	2.15	0.61
3:C:558:PHE:CZ	3:C:562:VAL:HG21	2.35	0.61
3:C:734:TYR:CE1	3:C:770:VAL:HG21	2.35	0.61
1:A:591:THR:HG22	1:A:595:TYR:CE2	2.37	0.60
1:A:264:LYS:HE2	1:A:294:PHE:CE2	2.37	0.60
1:A:602:SER:HG	1:A:620:TRP:CD1	2.19	0.60
3:C:521:VAL:CG1	3:C:578:TYR:CD2	2.84	0.60
3:C:702:LEU:C	3:C:704:LEU:H	2.08	0.60
1:A:603:LYS:HG3	3:C:564:GLU:CG	2.15	0.60
3:C:424:ILE:O	3:C:428:LEU:CG	2.41	0.59
1:A:350:PHE:CE2	1:A:384:LEU:CD1	2.85	0.59
3:C:521:VAL:CG1	3:C:578:TYR:HE2	2.13	0.59
3:C:639:LEU:HD23	3:C:643:ALA:HB3	1.84	0.59
3:C:595:ARG:O	3:C:599:LEU:HG	2.03	0.59
3:C:580:TYR:CD1	3:C:590:MET:CG	2.86	0.58
1:A:603:LYS:HG2	3:C:564:GLU:HG3	1.82	0.58
3:C:577:ARG:HA	3:C:580:TYR:CZ	2.38	0.58
3:C:636:LEU:HD21	3:C:692:LEU:O	2.02	0.58
3:C:517:VAL:HG11	3:C:578:TYR:CD2	2.38	0.58
3:C:435:GLN:O	3:C:439:ILE:HG23	2.03	0.58
3:C:622:ASN:HB3	3:C:625:THR:HG23	1.85	0.58
3:C:580:TYR:CD2	3:C:590:MET:CG	2.87	0.58
3:C:641:LYS:HG2	3:C:641:LYS:O	2.04	0.58
1:A:343:SER:CB	1:A:383:ILE:CB	2.82	0.57
1:A:602:SER:HB2	3:C:560:ALA:CB	2.20	0.57
1:A:343:SER:O	1:A:383:ILE:CG2	2.53	0.57
3:C:628:TRP:HZ3	3:C:689:PHE:HB2	1.68	0.57
1:A:351:GLN:HG2	1:A:391:LEU:HD11	1.85	0.56
1:A:47:GLU:OE1	1:A:135:LEU:CD2	2.53	0.56
3:C:546:ILE:HG13	3:C:550:LYS:HG3	1.86	0.56
1:A:350:PHE:HE1	1:A:359:VAL:CG2	2.05	0.56
1:A:80:SER:O	1:A:83:THR:HG22	2.06	0.56
3:C:638:TYR:CD2	3:C:643:ALA:HB2	2.41	0.55
3:C:742:GLU:O	3:C:746:ILE:HG13	2.06	0.55
1:A:351:GLN:CG	1:A:391:LEU:HD11	2.36	0.55
3:C:661:ARG:HG2	3:C:665:HIS:CD2	2.40	0.55
3:C:591:GLU:HB3	3:C:595:ARG:NH1	2.22	0.55
3:C:758:PRO:CA	3:C:761:ARG:HB2	2.32	0.55
3:C:726:LEU:HD23	3:C:759:THR:CG2	2.37	0.55
3:C:746:ILE:HG21	3:C:767:ILE:HG21	1.87	0.55
3:C:766:GLU:HG3	3:C:769:ARG:HH12	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:620:TRP:CE3	3:C:556:THR:CG2	2.90	0.55
1:A:243:PHE:CD1	1:A:280:LEU:HD11	2.42	0.54
1:A:83:THR:CG2	2:B:95:TYR:CZ	2.90	0.54
1:A:226:TYR:OH	1:A:255:PRO:CB	2.56	0.54
3:C:521:VAL:HG13	3:C:579:LEU:HG	1.89	0.54
3:C:437:ARG:O	3:C:441:ILE:HG13	2.08	0.54
3:C:448:GLU:HG2	3:C:519:VAL:HG12	1.89	0.54
3:C:702:LEU:C	3:C:704:LEU:N	2.65	0.54
3:C:714:LYS:HE2	3:C:740:LEU:O	2.07	0.54
3:C:742:GLU:O	3:C:746:ILE:N	2.34	0.54
1:A:606:HIS:CE1	3:C:442:ARG:HA	2.43	0.54
2:B:12:SER:O	2:B:16:ARG:HG3	2.08	0.54
1:A:343:SER:C	1:A:383:ILE:CG2	2.76	0.53
3:C:539:GLY:O	3:C:543:VAL:HB	2.08	0.53
3:C:569:TRP:CE2	3:C:600:ASN:CG	2.86	0.53
3:C:733:PHE:HE1	3:C:771:ARG:NE	2.07	0.53
3:C:746:ILE:HD11	3:C:771:ARG:HH12	1.72	0.53
3:C:449:TYR:CE1	3:C:562:VAL:HG22	2.42	0.53
3:C:556:THR:HA	3:C:592:LEU:HD11	1.91	0.53
3:C:758:PRO:O	3:C:761:ARG:HB2	2.09	0.53
3:C:617:GLN:NE2	3:C:620:LYS:HD2	2.21	0.52
3:C:569:TRP:HA	3:C:572:LEU:HB3	1.92	0.52
3:C:682:SER:HA	3:C:685:ASN:HD22	1.74	0.52
1:A:603:LYS:CG	3:C:564:GLU:CG	2.75	0.52
3:C:559:PHE:CE2	3:C:592:LEU:HB3	2.45	0.52
1:A:620:TRP:CZ3	3:C:556:THR:HG21	2.42	0.52
1:A:83:THR:HG23	1:A:84:ALA:N	2.24	0.52
1:A:347:GLN:HG2	1:A:387:SER:HB3	0.69	0.52
3:C:580:TYR:CD1	3:C:590:MET:SD	3.03	0.52
3:C:521:VAL:HG12	3:C:578:TYR:CE2	2.41	0.51
3:C:714:LYS:HB2	3:C:740:LEU:HD23	1.92	0.51
3:C:679:SER:HG	3:C:682:SER:H	1.56	0.51
3:C:657:VAL:CG2	3:C:705:ILE:HG21	2.40	0.50
3:C:455:VAL:HG11	3:C:517:VAL:HG22	1.93	0.50
3:C:639:LEU:HD23	3:C:643:ALA:CB	2.41	0.50
1:A:84:ALA:HA	2:B:95:TYR:CD2	2.46	0.50
1:A:351:GLN:HG2	1:A:391:LEU:CD1	2.42	0.50
3:C:526:LYS:O	3:C:530:GLU:HB2	2.12	0.50
1:A:343:SER:HB2	1:A:383:ILE:CB	2.35	0.49
3:C:562:VAL:HG12	3:C:572:LEU:HD11	1.94	0.49
3:C:738:PHE:CD1	3:C:738:PHE:N	2.80	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:449:TYR:HE1	3:C:562:VAL:HG22	1.77	0.49
3:C:679:SER:OG	3:C:682:SER:N	2.33	0.49
3:C:747:VAL:HA	3:C:750:ILE:HD12	1.94	0.49
1:A:376:THR:OG1	1:A:379:LEU:HG	2.12	0.48
3:C:747:VAL:HG22	3:C:764:ILE:HG23	1.95	0.48
3:C:431:GLU:O	3:C:434:MET:HB2	2.13	0.48
3:C:555:GLN:O	3:C:589:LEU:HD21	2.13	0.48
1:A:606:HIS:HB2	3:C:561:MET:SD	2.53	0.48
1:A:591:THR:O	1:A:595:TYR:CD2	2.66	0.48
1:A:350:PHE:CZ	1:A:384:LEU:HD12	2.45	0.48
1:A:83:THR:HG23	2:B:95:TYR:CD1	2.48	0.48
1:A:243:PHE:CD1	1:A:280:LEU:HD21	2.49	0.48
1:A:606:HIS:O	3:C:442:ARG:HG2	2.14	0.48
3:C:738:PHE:HD2	3:C:746:ILE:HD13	1.79	0.47
3:C:765:ASP:O	3:C:769:ARG:HB2	2.14	0.47
1:A:608:CYS:O	3:C:442:ARG:NH2	2.46	0.47
3:C:517:VAL:HB	3:C:521:VAL:HG21	1.95	0.47
3:C:620:LYS:HE2	3:C:675:TRP:CZ2	2.49	0.47
1:A:85:ILE:O	1:A:96:ARG:NH1	2.47	0.47
3:C:448:GLU:HB3	3:C:520:PHE:CB	2.41	0.47
3:C:681:LYS:O	3:C:681:LYS:HG3	2.14	0.47
3:C:569:TRP:O	3:C:572:LEU:HB3	2.15	0.46
1:A:591:THR:O	1:A:595:TYR:HD2	1.98	0.46
3:C:673:LEU:O	3:C:752:LEU:HD11	2.15	0.46
3:C:733:PHE:CG	3:C:767:ILE:HG12	2.50	0.46
3:C:589:LEU:HD23	3:C:589:LEU:HA	1.53	0.46
3:C:743:THR:HG21	3:C:768:ARG:HA	1.98	0.46
1:A:84:ALA:CA	2:B:95:TYR:HB2	2.45	0.46
3:C:690:ILE:O	3:C:693:LEU:CA	2.63	0.46
3:C:714:LYS:NZ	3:C:745:GLU:OE2	2.49	0.46
3:C:430:PHE:HE2	3:C:543:VAL:HG13	1.80	0.46
3:C:456:GLN:OE1	3:C:514:PRO:HA	2.15	0.46
3:C:758:PRO:C	3:C:761:ARG:HB2	2.41	0.46
1:A:243:PHE:HD1	1:A:280:LEU:HD21	1.81	0.46
3:C:635:LEU:O	3:C:638:TYR:HB3	2.16	0.45
3:C:683:SER:O	3:C:687:SER:N	2.47	0.45
3:C:702:LEU:O	3:C:704:LEU:N	2.49	0.45
1:A:83:THR:HG23	2:B:95:TYR:CE1	2.51	0.45
3:C:455:VAL:HG21	3:C:578:TYR:OH	2.17	0.45
3:C:628:TRP:CZ3	3:C:689:PHE:HB2	2.50	0.45
1:A:19:LYS:HE2	1:A:23:GLU:OE1	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:525:ALA:HB2	3:C:582:GLU:OE1	2.17	0.45
3:C:570:THR:HB	3:C:650:LEU:HD21	1.98	0.45
3:C:709:ARG:HG2	3:C:713:GLU:OE2	2.16	0.45
3:C:638:TYR:O	3:C:642:THR:HB	2.17	0.45
3:C:577:ARG:HA	3:C:580:TYR:CE2	2.52	0.44
1:A:83:THR:OG1	2:B:95:TYR:OH	2.08	0.44
3:C:448:GLU:HG2	3:C:519:VAL:CG1	2.48	0.44
3:C:684:GLU:HG2	3:C:723:THR:HG22	1.99	0.44
3:C:747:VAL:O	3:C:750:ILE:HB	2.18	0.44
3:C:749:TRP:HA	3:C:749:TRP:CE3	2.53	0.44
1:A:343:SER:O	1:A:383:ILE:HG22	2.17	0.44
3:C:627:ASP:OD1	3:C:629:ASP:HB2	2.18	0.44
3:C:449:TYR:HA	3:C:520:PHE:CD2	2.53	0.44
3:C:452:LEU:HA	3:C:518:SER:HA	1.98	0.44
3:C:455:VAL:CG1	3:C:517:VAL:HG22	2.48	0.44
1:A:1:MET:HE3	1:A:28:LEU:HD21	1.99	0.44
3:C:638:TYR:HA	3:C:642:THR:HB	2.00	0.44
3:C:747:VAL:HG11	3:C:768:ARG:NH1	2.33	0.44
1:A:603:LYS:HE3	3:C:564:GLU:OE1	2.18	0.43
2:B:16:ARG:HG2	2:B:76:LEU:HD13	1.99	0.43
3:C:663:ILE:O	3:C:667:LEU:HB3	2.18	0.43
3:C:515:ILE:HD11	3:C:574:LYS:HD2	2.00	0.43
3:C:569:TRP:O	3:C:569:TRP:CE3	2.72	0.43
3:C:636:LEU:HD12	3:C:639:LEU:HD12	2.00	0.43
3:C:559:PHE:CD2	3:C:592:LEU:HD13	2.53	0.43
3:C:562:VAL:HG12	3:C:572:LEU:CD1	2.49	0.43
3:C:734:TYR:OH	3:C:766:GLU:HG2	2.19	0.43
3:C:569:TRP:HH2	3:C:597:LEU:CD1	2.32	0.43
1:A:384:LEU:O	1:A:387:SER:OG	2.33	0.42
3:C:727:LYS:HE3	3:C:731:GLU:OE2	2.20	0.42
3:C:681:LYS:HA	3:C:681:LYS:HD2	1.75	0.42
3:C:689:PHE:O	3:C:692:LEU:HB2	2.20	0.42
3:C:609:TYR:CD1	3:C:609:TYR:C	2.97	0.42
3:C:743:THR:OG1	3:C:771:ARG:HG3	2.18	0.42
1:A:610:GLU:CD	3:C:442:ARG:HH12	2.27	0.42
2:B:42:LEU:HD23	2:B:80:VAL:CG2	2.49	0.42
3:C:558:PHE:CD2	3:C:558:PHE:C	2.98	0.42
3:C:690:ILE:C	3:C:693:LEU:H	2.27	0.42
1:A:319:MET:HE1	1:A:327:LEU:HD11	2.01	0.42
3:C:615:GLN:OE1	3:C:630:ILE:HB	2.20	0.42
3:C:573:TYR:CE1	3:C:593:ASN:ND2	2.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:579:LEU:HA	3:C:579:LEU:HD23	1.61	0.42
3:C:517:VAL:HG11	3:C:578:TYR:CG	2.55	0.41
3:C:622:ASN:O	3:C:623:GLU:C	2.63	0.41
1:A:343:SER:HB2	1:A:383:ILE:HG12	2.02	0.41
1:A:384:LEU:C	1:A:384:LEU:HD23	2.45	0.41
3:C:569:TRP:CE2	3:C:600:ASN:OD1	2.73	0.41
3:C:604:ILE:HG12	3:C:638:TYR:OH	2.20	0.41
3:C:728:ASP:O	3:C:731:GLU:HB2	2.20	0.41
1:A:313:LEU:HD23	1:A:316:MET:SD	2.60	0.41
1:A:347:GLN:HA	1:A:387:SER:CB	2.50	0.41
1:A:603:LYS:HG2	3:C:564:GLU:CB	2.51	0.41
3:C:672:ILE:HG22	3:C:749:TRP:HH2	1.85	0.41
3:C:615:GLN:HB2	3:C:630:ILE:CG2	2.50	0.41
3:C:733:PHE:HB2	3:C:738:PHE:CD2	2.55	0.41
3:C:747:VAL:HG11	3:C:768:ARG:HH11	1.86	0.41
1:A:83:THR:CG2	1:A:84:ALA:N	2.83	0.41
3:C:688:GLU:O	3:C:692:LEU:HG	2.21	0.41
1:A:87:GLU:O	2:B:98:TYR:HE1	1.98	0.41
3:C:574:LYS:HE3	3:C:574:LYS:HB2	1.82	0.41
2:B:16:ARG:HH22	2:B:76:LEU:H	1.67	0.41
3:C:698:GLU:HB2	3:C:712:ARG:NH2	2.35	0.41
3:C:654:THR:O	3:C:657:VAL:HB	2.20	0.40
3:C:733:PHE:CE2	3:C:767:ILE:HA	2.55	0.40
3:C:766:GLU:HA	3:C:769:ARG:HH12	1.86	0.40
3:C:538:ILE:HG12	3:C:542:LYS:HD2	2.04	0.40
3:C:694:LEU:HD23	3:C:694:LEU:HA	1.78	0.40
3:C:703:ASN:OD1	3:C:708:TYR:HE2	2.03	0.40
3:C:733:PHE:HE2	3:C:770:VAL:HB	1.85	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	621/699 (89%)	607 (98%)	14 (2%)	0	100	100
2	B	87/113 (77%)	87 (100%)	0	0	100	100
3	C	290/306 (95%)	273 (94%)	16 (6%)	1 (0%)	36	72
All	All	998/1118 (89%)	967 (97%)	30 (3%)	1 (0%)	48	83

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	538	ILE

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	584/645 (90%)	581 (100%)	3 (0%)	81	82
2	B	84/106 (79%)	84 (100%)	0	100	100
3	C	279/288 (97%)	266 (95%)	13 (5%)	23	44
All	All	947/1039 (91%)	931 (98%)	16 (2%)	53	68

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	VAL
1	A	256	LEU
1	A	602	SER
3	C	453	VAL
3	C	515	ILE
3	C	538	ILE
3	C	548	LEU
3	C	571	GLN
3	C	581	THR
3	C	591	GLU
3	C	601	LEU
3	C	660	ILE

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Mol	Chain	Res	Type
3	C	694	LEU
3	C	713	GLU
3	C	740	LEU
3	C	748	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	347	GLN
1	A	380	GLN
1	A	569	HIS
1	A	606	HIS
2	B	14	GLN
2	B	18	ASN
2	B	35	HIS
3	C	436	GLN
3	C	516	GLN
3	C	622	ASN
3	C	665	HIS
3	C	685	ASN

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

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	A	627/699 (89%)	1.37	160 (25%)  	174, 288, 375, 442	0
2	B	91/113 (80%)	0.88	17 (18%)  	164, 236, 316, 328	0
3	C	294/306 (96%)	0.60	25 (8%)  	170, 230, 306, 350	0
All	All	1012/1118 (90%)	1.10	202 (19%)  	164, 271, 365, 442	0

All (202) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	319	MET	8.6
1	A	230	PHE	8.3
1	A	222	TYR	6.0
1	A	72	LEU	5.9
1	A	3	PRO	5.8
1	A	254	THR	5.7
1	A	4	GLN	5.6
1	A	318	ALA	5.4
1	A	349	MET	5.3
1	A	362	VAL	5.1
1	A	86	SER	5.0
3	C	426	THR	5.0
1	A	604	LEU	4.8
1	A	240	ILE	4.8
1	A	190	VAL	4.6
1	A	597	GLN	4.6
1	A	193	MET	4.5
1	A	317	ASN	4.5
1	A	595	TYR	4.5
1	A	601	VAL	4.4
1	A	623	LEU	4.3
1	A	626	GLU	4.2
1	A	229	ILE	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	510	PHE	4.1
3	C	768	ARG	4.1
1	A	225	PHE	4.1
1	A	421	PHE	4.0
1	A	243	PHE	4.0
1	A	458	THR	4.0
1	A	37	TYR	4.0
1	A	406	VAL	3.8
2	B	41	LYS	3.8
2	B	75	ARG	3.8
1	A	482	SER	3.7
1	A	492	VAL	3.7
1	A	632	LEU	3.7
1	A	294	PHE	3.6
1	A	199	LYS	3.6
1	A	87	GLU	3.5
1	A	603	LYS	3.5
1	A	121	LEU	3.5
1	A	76	ILE	3.5
1	A	194	TRP	3.5
1	A	85	ILE	3.5
1	A	182	GLU	3.5
1	A	478	PRO	3.5
1	A	75	LEU	3.5
1	A	366	ILE	3.5
1	A	621	SER	3.4
1	A	148	ASN	3.4
2	B	23	LEU	3.4
2	B	35	HIS	3.3
2	B	34	TYR	3.3
1	A	149	SER	3.3
2	B	19	TYR	3.3
1	A	181	PHE	3.3
1	A	178	ILE	3.3
1	A	352	GLY	3.3
1	A	350	PHE	3.3
1	A	489	MET	3.2
1	A	261	ASN	3.2
1	A	363	VAL	3.2
1	A	66	MET	3.2
1	A	135	LEU	3.2
1	A	47	GLU	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	342	PHE	3.2
1	A	409	GLN	3.1
1	A	49	ASP	3.1
1	A	327	LEU	3.1
3	C	624	ARG	3.1
3	C	564	GLU	3.1
1	A	255	PRO	3.1
1	A	239	SER	3.1
1	A	300	ILE	3.1
1	A	417	PHE	3.1
1	A	315	PHE	3.1
1	A	437	ALA	3.1
1	A	7	LEU	3.0
1	A	496	LEU	3.0
1	A	628	SER	3.0
1	A	481	ILE	3.0
1	A	196	LEU	3.0
1	A	124	PHE	3.0
1	A	345	ILE	2.9
1	A	120	MET	2.9
1	A	627	LEU	2.9
2	B	31	ASP	2.9
1	A	441	LEU	2.9
1	A	191	SER	2.9
1	A	197	PRO	2.9
1	A	40	LEU	2.9
3	C	632	ILE	2.8
1	A	277	SER	2.8
1	A	476	PHE	2.8
1	A	599	GLY	2.8
1	A	15	ARG	2.8
1	A	479	ALA	2.8
1	A	145	LEU	2.8
1	A	307	ASN	2.8
1	A	351	GLN	2.8
1	A	625	LEU	2.8
3	C	611	LEU	2.8
1	A	258	THR	2.8
1	A	158	LEU	2.8
3	C	684	GLU	2.7
1	A	401	VAL	2.7
1	A	237	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
3	C	565	LYS	2.7
1	A	195	LYS	2.7
1	A	422	ASN	2.7
1	A	314	LYS	2.6
2	B	78	GLU	2.6
1	A	152	LYS	2.6
3	C	751	ILE	2.6
1	A	312	ASN	2.6
1	A	346	THR	2.6
1	A	438	ASN	2.6
1	A	296	ASP	2.6
1	A	280	LEU	2.6
2	B	99	GLN	2.6
1	A	44	MET	2.6
1	A	328	LYS	2.5
3	C	767	ILE	2.5
3	C	641	LYS	2.5
1	A	615	VAL	2.5
1	A	157	TRP	2.5
1	A	263	THR	2.5
2	B	67	ASP	2.5
1	A	608	CYS	2.5
1	A	390	LYS	2.5
1	A	115	ILE	2.4
1	A	246	TYR	2.4
1	A	533	TYR	2.4
3	C	678	ILE	2.4
1	A	42	CYS	2.4
1	A	291	ILE	2.4
3	C	764	ILE	2.4
3	C	578	TYR	2.4
1	A	224	PRO	2.4
3	C	774	ALA	2.4
2	B	26	PHE	2.4
1	A	43	VAL	2.4
1	A	288	GLU	2.4
2	B	11	LEU	2.4
3	C	687	SER	2.4
1	A	594	ILE	2.4
1	A	162	LEU	2.3
3	C	688	GLU	2.3
3	C	526	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	513	LEU	2.3
1	A	598	LEU	2.3
1	A	564	PRO	2.3
2	B	16	ARG	2.3
1	A	344	ILE	2.3
1	A	541	ILE	2.3
1	A	228	VAL	2.3
1	A	418	TRP	2.3
1	A	250	SER	2.3
1	A	580	ILE	2.3
1	A	78	GLU	2.3
1	A	316	MET	2.3
1	A	365	PHE	2.3
1	A	156	SER	2.3
3	C	587	LEU	2.3
1	A	434	MET	2.3
2	B	9	ASP	2.3
1	A	348	ASN	2.2
1	A	429	LEU	2.2
3	C	516	GLN	2.2
1	A	466	TYR	2.2
1	A	259	LYS	2.2
1	A	442	ASN	2.2
1	A	633	ILE	2.2
1	A	293	SER	2.2
1	A	468	LEU	2.2
1	A	73	VAL	2.2
1	A	306	LEU	2.2
3	C	548	LEU	2.2
1	A	126	ARG	2.1
1	A	292	LYS	2.1
1	A	407	LYS	2.1
1	A	518	THR	2.1
2	B	66	ILE	2.1
2	B	62	SER	2.1
1	A	610	GLU	2.1
1	A	341	TYR	2.1
1	A	396	ILE	2.1
1	A	147	ALA	2.1
3	C	707	THR	2.1
1	A	8	CYS	2.1
3	C	585	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
3	C	763	CYS	2.1
1	A	12	PHE	2.1
2	B	38	ILE	2.1
1	A	14	THR	2.1
1	A	448	GLU	2.0
1	A	624	GLU	2.0
1	A	90	VAL	2.0
3	C	573	TYR	2.0
1	A	265	TRP	2.0
1	A	69	SER	2.0
1	A	324	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.