



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:12 PM UTC

PDB ID : 1WIP / pdb_00001wip
Title : STRUCTURE OF T-CELL SURFACE GLYCOPROTEIN CD4, MONO-CLINIC CRYSTAL FORM
Authors : Wu, H.; Kwong, P.D.; Hendrickson, W.A.
Deposited on : 1996-12-18
Resolution : 4.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

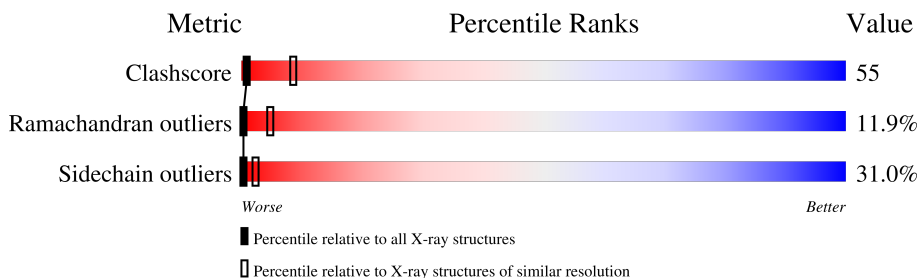
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	1129 (4.20-3.80)
Ramachandran outliers	187476	1064 (4.20-3.80)
Sidechain outliers	187428	1055 (4.20-3.80)

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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	363	20% 47% 27% 6%
1	B	363	21% 48% 25% 5%

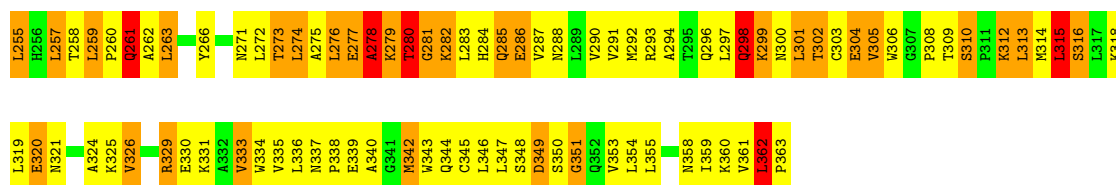
2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5624 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 T-CELL SURFACE GLYCOPROTEIN CD4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	363	Total	C	N	O	S	0	0	0
			2812	1784	479	539	10			
1	B	363	Total	C	N	O	S	0	0	0
			2812	1784	479	539	10			



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	100.50 Å 123.40 Å 100.60 Å 90.00° 103.40° 90.00°	Depositor
Resolution (Å)	8.00 – 4.00	Depositor
% Data completeness (in resolution range)	76.0 (8.00-4.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.452 , 0.427	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5624	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

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	0.89	5/2861 (0.2%)	1.45	48/3871 (1.2%)
1	B	0.87	4/2861 (0.1%)	1.36	40/3871 (1.0%)
All	All	0.88	9/5722 (0.2%)	1.40	88/7742 (1.1%)

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
1	A	0	1
1	B	0	1
All	All	0	2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	178	ALA	C-N	17.89	1.57	1.33
1	A	202	PHE	C-O	-8.28	1.13	1.24
1	B	202	PHE	C-O	-7.56	1.14	1.24
1	A	178	ALA	C-N	7.51	1.44	1.33
1	B	249	MET	SD-CE	-6.83	1.62	1.79
1	A	260	PRO	C-O	-6.72	1.15	1.24
1	A	178	ALA	CA-CB	-6.58	1.42	1.53
1	B	260	PRO	C-O	-5.36	1.17	1.23
1	A	207	LEU	CA-C	5.12	1.59	1.52

All (88) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	178	ALA	O-C-N	-27.61	85.88	122.59
1	B	208	THR	N-CA-C	12.49	133.62	113.89
1	A	208	THR	N-CA-C	11.81	131.33	113.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	178	ALA	O-C-N	-11.37	107.47	122.59
1	A	87	GLU	N-CA-C	8.73	129.40	110.80
1	B	87	GLU	N-CA-C	8.58	129.07	110.80
1	B	226	TRP	N-CA-C	8.53	119.16	107.73
1	B	301	LEU	N-CA-C	8.53	121.98	109.69
1	A	277	GLU	O-C-N	8.51	133.91	122.59
1	A	226	TRP	N-CA-C	8.46	119.06	107.73
1	B	277	GLU	O-C-N	8.41	133.78	122.59
1	B	67	PHE	CA-C-N	8.10	128.03	119.85
1	B	67	PHE	C-N-CA	8.10	128.03	119.85
1	B	362	LEU	N-CA-C	7.96	127.39	109.81
1	A	301	LEU	N-CA-C	7.71	121.06	109.95
1	A	202	PHE	N-CA-C	7.43	126.64	110.80
1	B	106	THR	N-CA-C	7.21	126.15	110.80
1	B	105	ASP	N-CA-C	7.15	126.02	110.80
1	B	110	GLN	N-CA-C	7.12	119.93	110.53
1	A	106	THR	N-CA-C	7.09	125.89	110.80
1	A	67	PHE	N-CA-C	-6.95	98.67	109.79
1	A	47	GLY	CA-C-N	6.75	126.37	119.56
1	A	47	GLY	C-N-CA	6.75	126.37	119.56
1	A	105	ASP	N-CA-C	6.71	125.08	110.80
1	A	67	PHE	CA-C-N	6.70	126.61	119.85
1	A	67	PHE	C-N-CA	6.70	126.61	119.85
1	B	239	LYS	N-CA-C	-6.63	103.98	111.07
1	B	67	PHE	N-CA-C	-6.55	99.31	109.79
1	B	142	LYS	N-CA-C	-6.55	105.27	113.20
1	A	278	ALA	O-C-N	6.50	131.23	122.59
1	A	142	LYS	N-CA-C	-6.47	105.49	113.19
1	B	278	ALA	O-C-N	6.45	131.17	122.59
1	B	315	LEU	N-CA-C	-6.45	99.31	109.50
1	A	242	THR	N-CA-C	-6.42	97.13	110.80
1	A	244	ASP	CA-C-N	-6.36	114.23	120.52
1	A	244	ASP	C-N-CA	-6.36	114.23	120.52
1	A	253	LEU	N-CA-C	6.23	123.57	109.81
1	A	251	LYS	N-CA-C	-6.22	105.51	113.23
1	B	244	ASP	CA-C-N	-6.17	114.38	120.98
1	B	244	ASP	C-N-CA	-6.17	114.38	120.98
1	B	261	GLN	N-CA-C	6.16	123.91	110.80
1	A	362	LEU	N-CA-C	6.14	123.38	109.81
1	A	203	THR	N-CA-C	6.14	123.88	110.80
1	B	21	LYS	N-CA-C	6.11	123.81	110.80
1	A	109	LEU	N-CA-C	-6.05	101.04	110.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	78	ASP	N-CA-C	-6.03	105.97	113.38
1	B	133	PRO	N-CA-C	-6.00	106.36	113.86
1	B	253	LEU	N-CA-C	5.98	123.02	109.81
1	B	280	THR	O-C-N	5.96	130.52	122.59
1	A	280	THR	O-C-N	5.96	130.52	122.59
1	A	133	PRO	N-CA-C	-5.77	106.64	113.86
1	A	184	SER	N-CA-C	-5.74	101.99	110.48
1	A	315	LEU	N-CA-C	-5.69	100.50	109.50
1	B	325	LYS	N-CA-C	-5.69	106.17	113.23
1	A	21	LYS	N-CA-C	5.68	122.91	110.80
1	B	121	PRO	CA-C-N	5.61	126.86	119.84
1	B	121	PRO	C-N-CA	5.61	126.86	119.84
1	A	263	LEU	CA-C-N	5.60	125.70	119.32
1	A	263	LEU	C-N-CA	5.60	125.70	119.32
1	A	208	THR	CB-CA-C	-5.59	100.26	109.99
1	B	298	GLN	N-CA-C	-5.59	98.90	110.80
1	B	340	ALA	N-CA-C	5.58	118.90	110.64
1	A	239	LYS	N-CA-C	-5.49	105.19	111.07
1	A	101	THR	N-CA-C	5.43	117.51	108.99
1	A	358	ASN	N-CA-C	5.38	118.45	110.48
1	B	193	GLN	N-CA-CB	5.35	117.94	109.71
1	A	121	PRO	CA-C-N	5.31	126.48	119.84
1	A	121	PRO	C-N-CA	5.31	126.48	119.84
1	B	101	THR	N-CA-C	5.29	117.30	108.99
1	B	251	LYS	N-CA-C	-5.23	106.74	113.23
1	B	47	GLY	CA-C-N	5.23	124.84	119.56
1	B	47	GLY	C-N-CA	5.23	124.84	119.56
1	A	123	GLY	N-CA-C	5.20	125.49	113.18
1	A	252	LYS	N-CA-C	5.18	117.94	109.59
1	A	324	ALA	N-CA-C	5.17	121.81	110.80
1	B	304	GLU	N-CA-C	5.15	117.07	108.99
1	A	17	THR	N-CA-C	-5.14	101.56	108.19
1	A	110	GLN	N-CA-C	5.13	116.90	110.24
1	B	260	PRO	N-CA-C	-5.09	103.20	111.14
1	B	184	SER	N-CA-C	-5.06	102.27	110.32
1	A	281	GLY	N-CA-C	-5.06	104.36	111.09
1	A	361	VAL	CA-C-O	-5.05	114.66	121.32
1	B	263	LEU	CA-C-N	5.05	125.07	119.32
1	B	263	LEU	C-N-CA	5.05	125.07	119.32
1	A	110	GLN	N-CA-CB	-5.03	101.86	109.92
1	B	281	GLY	N-CA-C	-5.03	104.40	111.09
1	A	135	GLY	N-CA-C	5.02	120.11	111.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	353	VAL	N-CA-C	5.02	115.70	108.93

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	178	ALA	Mainchain
1	B	178	ALA	Mainchain

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	2812	0	2873	307	19
1	B	2812	0	2873	332	19
All	All	5624	0	5746	627	19

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 55.

All (627) 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:112:GLN:OE1	1:B:281:GLY:CA	1.68	1.41
1:A:109:LEU:HD23	1:A:283:LEU:CD2	1.67	1.22
1:B:112:GLN:OE1	1:B:281:GLY:HA3	1.44	1.14
1:B:276:LEU:HD23	1:B:276:LEU:N	1.63	1.11
1:A:276:LEU:H	1:A:276:LEU:CD2	1.63	1.09
1:B:112:GLN:OE1	1:B:281:GLY:HA2	1.32	1.08
1:A:245:PRO:HB3	1:A:266:TYR:HE2	1.09	1.08
1:B:104:SER:HA	1:B:279:LYS:NZ	1.67	1.08
1:B:276:LEU:H	1:B:276:LEU:CD2	1.63	1.08
1:A:276:LEU:HD23	1:A:276:LEU:N	1.63	1.07
1:B:178:ALA:HA	1:B:283:LEU:HD22	1.20	1.07
1:A:234:LYS:HA	1:A:253:LEU:HG	1.35	1.05

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:314:MET:SD	1:B:329:ARG:HG3	1.96	1.05
1:A:318:LYS:HE3	1:A:320:GLU:HA	1.41	1.02
1:B:234:LYS:HA	1:B:253:LEU:HG	1.39	1.02
1:B:318:LYS:HE3	1:B:320:GLU:HA	1.39	1.02
1:A:276:LEU:H	1:A:276:LEU:HD23	0.84	1.00
1:B:276:LEU:HD23	1:B:276:LEU:H	0.84	1.00
1:B:177:LEU:HD22	1:B:283:LEU:HD11	1.41	0.98
1:A:245:PRO:HB3	1:A:266:TYR:CE2	1.99	0.98
1:A:178:ALA:C	1:A:179:PHE:O	2.06	0.98
1:A:276:LEU:CD2	1:A:276:LEU:N	2.19	0.98
1:A:109:LEU:HD23	1:A:283:LEU:HD21	1.45	0.96
1:A:314:MET:SD	1:A:329:ARG:HG3	2.05	0.96
1:B:276:LEU:N	1:B:276:LEU:CD2	2.19	0.94
1:A:309:THR:HB	1:A:313:LEU:HD11	1.46	0.94
1:B:104:SER:HA	1:B:279:LYS:HZ2	1.32	0.93
1:B:178:ALA:HA	1:B:283:LEU:CD2	1.98	0.93
1:A:218:GLU:HG2	1:A:219:ARG:HD2	1.51	0.92
1:A:344:GLN:HE22	1:B:344:GLN:HE22	1.19	0.89
1:B:178:ALA:CA	1:B:283:LEU:HD22	2.03	0.89
1:B:245:PRO:HG3	1:B:266:TYR:HE2	1.39	0.86
1:A:54:ARG:HD3	1:A:73:ASN:HB2	1.56	0.85
1:A:37:LEU:HA	1:A:46:LYS:HA	1.58	0.85
1:B:263:LEU:HB2	1:B:266:TYR:CE1	2.11	0.85
1:A:245:PRO:CB	1:A:266:TYR:HE2	1.89	0.85
1:B:124:SER:HB3	1:B:163:GLN:HE22	1.42	0.84
1:B:109:LEU:HD21	1:B:276:LEU:HD23	1.58	0.83
1:B:218:GLU:HG2	1:B:219:ARG:HD2	1.59	0.83
1:B:109:LEU:HD21	1:B:276:LEU:CD2	2.09	0.83
1:B:37:LEU:HA	1:B:46:LYS:HA	1.59	0.83
1:A:178:ALA:O	1:A:179:PHE:O	1.94	0.82
1:B:313:LEU:HB3	1:B:348:SER:O	1.79	0.82
1:B:319:LEU:CD1	1:B:321:ASN:HB2	2.10	0.81
1:A:111:GLY:O	1:A:148:GLN:HA	1.80	0.81
1:B:107:HIS:CE1	1:B:205:GLU:OE1	2.34	0.81
1:A:213:LEU:HB2	1:A:272:LEU:HD12	1.63	0.81
1:A:315:LEU:HD21	1:A:345:CYS:SG	2.21	0.81
1:A:244:ASP:HB3	1:A:245:PRO:CD	2.12	0.80
1:B:303:CYS:HG	1:B:345:CYS:HG	1.25	0.80
1:A:2:LYS:HB2	1:A:93:VAL:HG13	1.62	0.80
1:B:130:CYS:HG	1:B:159:CYS:HG	1.27	0.80
1:A:224:LYS:HA	1:A:226:TRP:CH2	2.17	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:ARG:O	1:A:62:TRP:HB2	1.82	0.79
1:B:130:CYS:SG	1:B:144:LEU:HD11	2.23	0.79
1:A:110:GLN:HE22	1:A:178:ALA:HB1	1.47	0.79
1:B:309:THR:HB	1:B:313:LEU:HD11	1.64	0.79
1:A:245:PRO:HG2	1:A:247:LEU:HD21	1.63	0.79
1:B:33:GLN:HA	1:B:33:GLN:OE1	1.82	0.79
1:B:54:ARG:HD3	1:B:73:ASN:HB2	1.64	0.78
1:A:244:ASP:HB3	1:A:245:PRO:HD2	1.66	0.78
1:A:109:LEU:HD23	1:A:283:LEU:CG	2.12	0.78
1:A:109:LEU:HD23	1:A:283:LEU:HD23	1.62	0.77
1:A:344:GLN:HE22	1:B:344:GLN:NE2	1.82	0.77
1:B:244:ASP:HB3	1:B:245:PRO:HD3	1.66	0.77
1:B:160:THR:HG21	1:B:167:LYS:HE3	1.67	0.77
1:B:245:PRO:HG3	1:B:266:TYR:CE2	2.20	0.77
1:B:59:ARG:O	1:B:62:TRP:HB2	1.86	0.76
1:A:76:ILE:HD12	1:A:76:ILE:H	1.49	0.76
1:A:29:LYS:HD3	1:A:33:GLN:OE1	1.87	0.75
1:A:178:ALA:O	1:A:179:PHE:C	2.21	0.75
1:B:245:PRO:HG2	1:B:247:LEU:HD21	1.67	0.75
1:A:162:LEU:HD23	1:A:163:GLN:N	2.02	0.75
1:B:76:ILE:HD12	1:B:76:ILE:H	1.50	0.75
1:B:277:GLU:HG3	1:B:278:ALA:N	2.02	0.74
1:B:213:LEU:HB2	1:B:272:LEU:HD12	1.66	0.74
1:A:277:GLU:HG3	1:A:278:ALA:N	2.02	0.74
1:B:162:LEU:HD23	1:B:163:GLN:N	2.01	0.74
1:A:160:THR:HG21	1:A:167:LYS:HE3	1.68	0.74
1:A:185:ILE:HG12	1:A:288:ASN:HB2	1.70	0.74
1:A:344:GLN:NE2	1:B:344:GLN:HE22	1.86	0.74
1:B:224:LYS:HA	1:B:226:TRP:CH2	2.23	0.74
1:A:292:MET:HE2	1:A:345:CYS:SG	2.28	0.73
1:B:314:MET:SD	1:B:329:ARG:CG	2.75	0.73
1:A:124:SER:HB2	1:A:163:GLN:HE22	1.52	0.73
1:B:214:TRP:CE3	1:B:226:TRP:CZ3	2.76	0.73
1:A:110:GLN:NE2	1:A:178:ALA:HB1	2.03	0.73
1:B:109:LEU:CD2	1:B:276:LEU:CD2	2.68	0.71
1:B:112:GLN:NE2	1:B:280:THR:O	2.22	0.71
1:B:299:LYS:H	1:B:299:LYS:HD3	1.55	0.71
1:A:36:ILE:O	1:A:46:LYS:HG3	1.91	0.71
1:B:315:LEU:HD21	1:B:345:CYS:SG	2.30	0.71
1:A:294:ALA:HA	1:A:303:CYS:SG	2.31	0.71
1:B:204:VAL:HG12	1:B:205:GLU:H	1.54	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:GLN:HG2	1:B:90:LYS:N	2.05	0.70
1:B:245:PRO:CG	1:B:266:TYR:HE2	2.05	0.70
1:A:310:SER:OG	1:A:312:LYS:HB2	1.92	0.70
1:B:107:HIS:HE1	1:B:205:GLU:OE1	1.72	0.70
1:B:245:PRO:HB3	1:B:266:TYR:CE2	2.26	0.70
1:B:1:LYS:HD2	1:B:92:GLU:HB2	1.74	0.69
1:A:187:TYR:CE2	1:A:354:LEU:HD22	2.27	0.69
1:B:107:HIS:ND1	1:B:278:ALA:CB	2.55	0.69
1:B:261:GLN:HE21	1:B:263:LEU:HD11	1.56	0.69
1:B:213:LEU:HD12	1:B:271:ASN:O	1.92	0.69
1:A:214:TRP:CE3	1:A:226:TRP:CZ3	2.80	0.69
1:A:130:CYS:HA	1:A:159:CYS:HA	1.75	0.69
1:A:193:GLN:HA	1:A:260:PRO:O	1.91	0.69
1:B:263:LEU:HB2	1:B:266:TYR:CD1	2.28	0.69
1:B:187:TYR:CE2	1:B:354:LEU:HD22	2.28	0.68
1:A:313:LEU:HB3	1:A:348:SER:O	1.94	0.68
1:B:290:VAL:HG13	1:B:306:TRP:O	1.92	0.68
1:A:177:LEU:HD23	1:A:178:ALA:N	2.09	0.68
1:A:210:SER:HB3	1:A:230:ASP:HA	1.73	0.68
1:A:17:THR:HG22	1:A:18:ALA:H	1.59	0.68
1:A:85:GLU:HG2	1:A:90:LYS:HG2	1.74	0.68
1:A:224:LYS:HA	1:A:226:TRP:CZ2	2.29	0.68
1:B:104:SER:HA	1:B:279:LYS:HZ3	1.54	0.68
1:A:89:GLN:HG2	1:A:90:LYS:N	2.09	0.68
1:A:110:GLN:NE2	1:A:178:ALA:CB	2.57	0.67
1:B:35:LYS:O	1:B:47:GLY:HA3	1.94	0.67
1:B:79:SER:HA	1:B:95:LEU:O	1.95	0.67
1:A:177:LEU:HD23	1:A:178:ALA:H	1.59	0.67
1:B:319:LEU:HD12	1:B:321:ASN:HB2	1.77	0.67
1:A:344:GLN:NE2	1:B:344:GLN:NE2	2.42	0.67
1:B:205:GLU:CG	1:B:208:THR:HG21	2.25	0.67
1:A:1:LYS:HD2	1:A:92:GLU:HB2	1.76	0.67
1:B:113:SER:HA	1:B:147:SER:O	1.95	0.66
1:B:111:GLY:O	1:B:148:GLN:HA	1.95	0.66
1:B:205:GLU:HG2	1:B:208:THR:HG21	1.76	0.66
1:A:27:HIS:CD2	1:A:85:GLU:HB2	2.31	0.66
1:A:223:SER:O	1:A:224:LYS:HB2	1.95	0.66
1:A:109:LEU:CD2	1:A:283:LEU:CD2	2.62	0.66
1:A:294:ALA:CA	1:A:303:CYS:SG	2.84	0.66
1:A:245:PRO:HG2	1:A:247:LEU:CD2	2.26	0.65
1:A:309:THR:HB	1:A:313:LEU:CD1	2.24	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:CYS:HG	1:A:159:CYS:HG	0.68	0.65
1:A:204:VAL:HG12	1:A:205:GLU:H	1.61	0.65
1:B:109:LEU:HD23	1:B:276:LEU:HD21	1.78	0.65
1:A:2:LYS:HG3	1:A:92:GLU:O	1.98	0.64
1:B:43:PHE:CD1	1:B:43:PHE:N	2.64	0.64
1:B:109:LEU:CD2	1:B:276:LEU:HD21	2.27	0.64
1:A:116:LEU:HB2	1:A:144:LEU:O	1.97	0.64
1:B:21:LYS:O	1:B:21:LYS:HG2	1.97	0.64
1:A:208:THR:O	1:A:276:LEU:CB	2.46	0.63
1:A:286:GLU:HG3	1:A:287:VAL:N	2.13	0.63
1:B:185:ILE:HG13	1:B:288:ASN:HB2	1.78	0.63
1:A:305:VAL:O	1:A:331:LYS:HE3	1.97	0.63
1:A:314:MET:SD	1:A:329:ARG:HD2	2.39	0.63
1:A:318:LYS:HG3	1:A:320:GLU:H	1.62	0.63
1:A:162:LEU:HD23	1:A:162:LEU:C	2.24	0.63
1:A:319:LEU:CD1	1:A:321:ASN:HB2	2.29	0.63
1:A:43:PHE:CD1	1:A:43:PHE:N	2.63	0.63
1:A:261:GLN:HE21	1:A:263:LEU:HD11	1.63	0.63
1:B:245:PRO:HB3	1:B:266:TYR:HE2	1.64	0.62
1:A:229:PHE:HE2	1:A:255:LEU:HD21	1.63	0.62
1:B:208:THR:O	1:B:276:LEU:CB	2.48	0.62
1:A:98:PHE:HZ	1:A:163:GLN:HE21	1.46	0.62
1:B:83:ILE:HA	1:B:91:GLU:O	1.99	0.62
1:B:210:SER:HB3	1:B:230:ASP:HA	1.82	0.62
1:B:310:SER:OG	1:B:312:LYS:HB2	1.99	0.62
1:B:362:LEU:HB2	1:B:363:PRO:HD3	1.82	0.62
1:A:59:ARG:HG2	1:A:62:TRP:CZ3	2.34	0.62
1:A:130:CYS:HG	1:A:159:CYS:CB	2.12	0.62
1:B:305:VAL:O	1:B:331:LYS:HE3	2.00	0.62
1:A:113:SER:HA	1:A:147:SER:O	1.99	0.62
1:A:305:VAL:CG1	1:A:309:THR:HG21	2.30	0.62
1:A:314:MET:SD	1:A:329:ARG:CG	2.83	0.61
1:B:224:LYS:HA	1:B:226:TRP:CZ2	2.35	0.61
1:B:318:LYS:HG3	1:B:320:GLU:H	1.64	0.61
1:A:196:PHE:HE2	1:A:259:LEU:HD12	1.63	0.61
1:A:265:GLN:HG2	1:A:266:TYR:CE1	2.35	0.61
1:B:263:LEU:HB2	1:B:266:TYR:HE1	1.65	0.61
1:A:338:PRO:HB2	1:A:361:VAL:HG21	1.81	0.61
1:A:347:LEU:HB3	1:A:355:LEU:HB2	1.81	0.61
1:B:36:ILE:O	1:B:46:LYS:HG3	2.00	0.61
1:B:106:THR:HG22	1:B:107:HIS:ND1	2.16	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:223:SER:O	1:B:224:LYS:HB2	2.01	0.61
1:A:28:TRP:NE1	1:A:69:LEU:HB2	2.16	0.60
1:B:124:SER:HB3	1:B:163:GLN:NE2	2.15	0.60
1:B:244:ASP:CB	1:B:245:PRO:HD3	2.31	0.60
1:B:121:PRO:HG2	1:B:124:SER:HB2	1.81	0.60
1:A:177:LEU:O	1:A:178:ALA:HB2	2.00	0.60
1:A:265:GLN:HG2	1:A:266:TYR:CD1	2.36	0.60
1:B:109:LEU:HD13	1:B:280:THR:C	2.27	0.60
1:B:43:PHE:N	1:B:43:PHE:HD1	2.00	0.60
1:B:314:MET:SD	1:B:329:ARG:HD2	2.41	0.60
1:A:43:PHE:N	1:A:43:PHE:HD1	1.98	0.60
1:A:83:ILE:HA	1:A:91:GLU:O	2.01	0.60
1:B:24:ILE:HD12	1:B:87:GLU:HB2	1.83	0.60
1:A:114:LEU:O	1:A:145:SER:HB3	2.02	0.60
1:A:106:THR:HG22	1:A:107:HIS:ND1	2.16	0.59
1:B:130:CYS:HA	1:B:159:CYS:HA	1.82	0.59
1:A:21:LYS:HG2	1:A:21:LYS:O	2.02	0.59
1:A:180:GLN:O	1:A:180:GLN:HG3	2.01	0.59
1:B:177:LEU:CD1	1:B:276:LEU:HD11	2.31	0.59
1:B:286:GLU:HG3	1:B:287:VAL:N	2.16	0.59
1:A:28:TRP:CE2	1:A:69:LEU:HB2	2.37	0.59
1:A:234:LYS:HA	1:A:253:LEU:CG	2.22	0.59
1:B:180:GLN:HB3	1:B:199:PRO:HG2	1.85	0.59
1:A:194:VAL:HB	1:A:259:LEU:HB3	1.83	0.59
1:A:229:PHE:CE2	1:A:255:LEU:HD21	2.38	0.59
1:A:111:GLY:HA2	1:A:148:GLN:CB	2.33	0.59
1:B:245:PRO:CB	1:B:266:TYR:HE2	2.16	0.59
1:B:347:LEU:O	1:B:354:LEU:HB2	2.03	0.58
1:A:7:LYS:HB2	1:A:10:ASP:HB2	1.84	0.58
1:B:245:PRO:HG2	1:B:247:LEU:CD2	2.32	0.58
1:A:37:LEU:HB2	1:A:45:THR:O	2.01	0.58
1:A:244:ASP:CB	1:A:245:PRO:HD2	2.34	0.58
1:A:109:LEU:HD23	1:A:283:LEU:HG	1.84	0.58
1:B:2:LYS:HD3	1:B:93:VAL:HG22	1.84	0.58
1:B:109:LEU:HD13	1:B:281:GLY:N	2.18	0.58
1:B:220:ALA:C	1:B:222:SER:H	2.12	0.58
1:B:17:THR:HG22	1:B:18:ALA:H	1.68	0.58
1:B:85:GLU:HG2	1:B:90:LYS:HG2	1.85	0.58
1:B:229:PHE:HE2	1:B:255:LEU:HD21	1.67	0.58
1:A:319:LEU:HB3	1:A:343:TRP:CE2	2.38	0.58
1:A:358:ASN:HD21	1:B:358:ASN:HD21	1.52	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:LYS:HB2	1:B:10:ASP:HB2	1.85	0.58
1:B:347:LEU:HB3	1:B:355:LEU:HB2	1.86	0.58
1:B:109:LEU:CD2	1:B:276:LEU:HD23	2.30	0.57
1:A:89:GLN:HG2	1:A:90:LYS:H	1.69	0.57
1:A:315:LEU:HA	1:A:346:LEU:O	2.04	0.57
1:A:342:MET:SD	1:A:360:LYS:HG2	2.45	0.57
1:B:116:LEU:HB2	1:B:144:LEU:O	2.03	0.57
1:B:194:VAL:HB	1:B:259:LEU:HB3	1.84	0.57
1:B:204:VAL:O	1:B:205:GLU:HB3	2.04	0.57
1:B:231:LEU:HG	1:B:236:VAL:HG13	1.86	0.57
1:B:241:VAL:HG12	1:B:242:THR:H	1.68	0.57
1:B:305:VAL:HG12	1:B:309:THR:HG21	1.85	0.57
1:B:296:GLN:HG2	1:B:298:GLN:O	2.05	0.57
1:A:14:LEU:HB2	1:A:69:LEU:HB3	1.87	0.56
1:A:79:SER:HA	1:A:95:LEU:O	2.05	0.56
1:B:28:TRP:NE1	1:B:69:LEU:HB2	2.20	0.56
1:A:273:THR:HG23	1:A:284:HIS:CD2	2.40	0.56
1:B:104:SER:CA	1:B:279:LYS:NZ	2.57	0.56
1:B:109:LEU:HD13	1:B:281:GLY:CA	2.35	0.56
1:B:204:VAL:O	1:B:205:GLU:CB	2.54	0.56
1:A:319:LEU:HD13	1:A:321:ASN:HB2	1.88	0.56
1:B:205:GLU:HG3	1:B:208:THR:CG2	2.35	0.56
1:A:106:THR:HG22	1:A:107:HIS:HD1	1.70	0.56
1:A:109:LEU:CD2	1:A:283:LEU:HD21	2.29	0.56
1:A:62:TRP:HA	1:A:66:ASN:O	2.05	0.56
1:B:279:LYS:O	1:B:280:THR:O	2.24	0.56
1:A:266:TYR:CD1	1:A:266:TYR:N	2.74	0.56
1:B:29:LYS:HA	1:B:34:ILE:O	2.06	0.56
1:A:196:PHE:CE2	1:A:259:LEU:HD12	2.41	0.55
1:A:279:LYS:O	1:A:280:THR:O	2.24	0.55
1:B:294:ALA:HA	1:B:303:CYS:SG	2.47	0.55
1:B:17:THR:O	1:B:18:ALA:HB2	2.06	0.55
1:B:177:LEU:HD23	1:B:178:ALA:N	2.21	0.55
1:B:273:THR:HG23	1:B:284:HIS:CD2	2.41	0.55
1:B:109:LEU:HD22	1:B:281:GLY:C	2.31	0.55
1:B:86:VAL:HG12	1:B:87:GLU:N	2.22	0.55
1:B:111:GLY:HA2	1:B:148:GLN:HB3	1.89	0.55
1:A:251:LYS:HB2	1:A:251:LYS:NZ	2.21	0.55
1:B:212:GLU:HB3	1:B:228:THR:HG23	1.88	0.55
1:A:296:GLN:HG2	1:A:298:GLN:O	2.06	0.55
1:B:112:GLN:HE22	1:B:280:THR:C	2.16	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:TRP:CD1	1:A:174:ILE:HD12	2.43	0.54
1:A:274:LEU:O	1:A:283:LEU:N	2.39	0.54
1:A:124:SER:CB	1:A:163:GLN:HE22	2.19	0.54
1:B:177:LEU:HD13	1:B:276:LEU:HD11	1.89	0.54
1:B:315:LEU:HA	1:B:346:LEU:O	2.06	0.54
1:B:37:LEU:HB2	1:B:45:THR:O	2.06	0.54
1:A:27:HIS:HD2	1:A:85:GLU:HB2	1.73	0.54
1:B:274:LEU:O	1:B:283:LEU:N	2.41	0.54
1:B:277:GLU:CG	1:B:278:ALA:N	2.59	0.54
1:A:259:LEU:O	1:A:260:PRO:O	2.26	0.54
1:A:298:GLN:OE1	1:A:299:LYS:HG2	2.08	0.54
1:A:305:VAL:HG12	1:A:309:THR:HG21	1.89	0.54
1:B:233:ASN:O	1:B:253:LEU:HD21	2.07	0.54
1:A:244:ASP:CB	1:A:245:PRO:CD	2.84	0.53
1:B:98:PHE:HZ	1:B:163:GLN:HE21	1.54	0.53
1:B:185:ILE:CG1	1:B:288:ASN:HB2	2.38	0.53
1:B:110:GLN:NE2	1:B:178:ALA:CB	2.71	0.53
1:A:19:SER:C	1:A:21:LYS:H	2.16	0.53
1:B:62:TRP:HA	1:B:66:ASN:O	2.07	0.53
1:B:100:LEU:HA	1:B:118:LEU:HD12	1.91	0.53
1:B:305:VAL:CG1	1:B:309:THR:HG21	2.37	0.53
1:A:141:GLY:C	1:A:143:THR:H	2.15	0.53
1:A:246:LYS:HB3	1:A:260:PRO:HG2	1.91	0.53
1:A:326:VAL:O	1:A:326:VAL:HG13	2.09	0.53
1:B:109:LEU:HD22	1:B:281:GLY:O	2.08	0.53
1:A:16:CYS:SG	1:A:84:CYS:SG	3.02	0.53
1:A:299:LYS:H	1:A:299:LYS:HD3	1.72	0.52
1:B:28:TRP:CE2	1:B:69:LEU:HB2	2.43	0.52
1:B:177:LEU:O	1:B:178:ALA:HB2	2.09	0.52
1:A:5:LEU:HD11	1:A:166:LYS:HB3	1.91	0.52
1:A:194:VAL:HB	1:A:259:LEU:CB	2.39	0.52
1:A:121:PRO:HG2	1:A:124:SER:OG	2.09	0.52
1:A:212:GLU:HB3	1:A:228:THR:HG23	1.92	0.52
1:A:358:ASN:ND2	1:B:358:ASN:ND2	2.58	0.52
1:B:157:TRP:CD1	1:B:174:ILE:HD12	2.45	0.52
1:A:52:ASN:OD1	1:A:53:ASP:N	2.43	0.52
1:A:87:GLU:O	1:A:88:ASP:CG	2.53	0.52
1:B:310:SER:C	1:B:312:LYS:H	2.18	0.52
1:A:164:ASN:C	1:A:166:LYS:H	2.16	0.52
1:B:2:LYS:HB2	1:B:93:VAL:HG13	1.91	0.52
1:B:177:LEU:HD23	1:B:178:ALA:H	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:276:LEU:N	1:B:276:LEU:HD22	2.21	0.52
1:B:292:MET:HE2	1:B:345:CYS:SG	2.49	0.52
1:B:345:CYS:C	1:B:346:LEU:HD12	2.35	0.52
1:A:80:ASP:O	1:A:94:GLN:HA	2.10	0.52
1:B:13:GLU:OE1	1:B:58:ARG:NH2	2.43	0.52
1:A:29:LYS:HA	1:A:34:ILE:O	2.11	0.51
1:B:19:SER:C	1:B:21:LYS:H	2.16	0.51
1:B:59:ARG:HH11	1:B:59:ARG:HB2	1.75	0.51
1:B:110:GLN:NE2	1:B:178:ALA:HB1	2.25	0.51
1:B:114:LEU:O	1:B:145:SER:HB3	2.11	0.51
1:A:161:VAL:HB	1:A:168:VAL:HG13	1.92	0.51
1:A:220:ALA:C	1:A:222:SER:H	2.17	0.51
1:A:233:ASN:O	1:A:253:LEU:HD21	2.11	0.51
1:B:27:HIS:CD2	1:B:85:GLU:HB2	2.46	0.51
1:B:206:LYS:HB3	1:B:207:LEU:HG	1.93	0.51
1:A:33:GLN:OE1	1:A:33:GLN:HA	2.09	0.51
1:B:244:ASP:CB	1:B:245:PRO:CD	2.88	0.51
1:A:266:TYR:O	1:A:289:LEU:HD23	2.11	0.51
1:A:291:VAL:O	1:A:305:VAL:HA	2.10	0.51
1:A:310:SER:O	1:A:313:LEU:HG	2.11	0.51
1:A:59:ARG:HA	1:A:62:TRP:CD2	2.45	0.51
1:B:109:LEU:HD13	1:B:281:GLY:HA3	1.92	0.51
1:A:109:LEU:CD2	1:A:283:LEU:HD23	2.37	0.51
1:A:315:LEU:HD13	1:A:333:VAL:HB	1.93	0.51
1:B:229:PHE:CE2	1:B:255:LEU:HD21	2.46	0.51
1:B:129:GLN:HA	1:B:139:GLN:HB3	1.93	0.51
1:A:204:VAL:O	1:A:205:GLU:CB	2.58	0.51
1:B:161:VAL:HB	1:B:168:VAL:HG13	1.92	0.51
1:A:212:GLU:HG3	1:A:214:TRP:HE1	1.76	0.50
1:A:342:MET:SD	1:A:360:LYS:HD2	2.50	0.50
1:A:132:SER:OG	1:A:135:GLY:O	2.28	0.50
1:A:26:PHE:CD2	1:A:67:PHE:HB3	2.46	0.50
1:A:86:VAL:HG12	1:A:87:GLU:N	2.26	0.50
1:A:100:LEU:HD12	1:A:118:LEU:HD12	1.92	0.50
1:A:206:LYS:O	1:A:208:THR:N	2.44	0.50
1:A:310:SER:C	1:A:312:LYS:H	2.19	0.50
1:B:319:LEU:HB3	1:B:343:TRP:CE2	2.46	0.50
1:A:219:ARG:NH1	1:A:306:TRP:HZ3	2.09	0.50
1:A:290:VAL:HG13	1:A:306:TRP:O	2.11	0.50
1:B:151:LEU:O	1:B:154:SER:HB2	2.10	0.50
1:B:182:ALA:O	1:B:286:GLU:HB3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:GLN:HG2	1:B:90:LYS:H	1.76	0.50
1:A:208:THR:HG22	1:A:277:GLU:H	1.77	0.50
1:A:345:CYS:C	1:A:346:LEU:HD12	2.36	0.50
1:B:106:THR:O	1:B:279:LYS:HB2	2.12	0.50
1:A:221:SER:O	1:A:223:SER:N	2.44	0.50
1:A:35:LYS:O	1:A:47:GLY:HA3	2.10	0.50
1:A:100:LEU:HD21	1:A:159:CYS:SG	2.52	0.50
1:A:24:ILE:HD12	1:A:87:GLU:HB2	1.93	0.49
1:A:12:VAL:HG21	1:A:95:LEU:HD11	1.94	0.49
1:A:130:CYS:N	1:A:138:ILE:O	2.45	0.49
1:A:326:VAL:HG11	1:A:335:VAL:HG22	1.93	0.49
1:B:294:ALA:CA	1:B:303:CYS:SG	3.00	0.49
1:B:14:LEU:HB2	1:B:69:LEU:HB3	1.95	0.49
1:A:109:LEU:HB3	1:A:283:LEU:HD23	1.93	0.49
1:A:46:LYS:HD2	1:A:55:ALA:HB3	1.94	0.49
1:A:54:ARG:HG2	1:A:72:LYS:HB2	1.94	0.49
1:A:44:LEU:HD13	1:A:62:TRP:HZ2	1.77	0.49
1:B:303:CYS:CB	1:B:345:CYS:HG	2.25	0.49
1:A:19:SER:O	1:A:21:LYS:N	2.38	0.49
1:A:204:VAL:O	1:A:205:GLU:HB3	2.11	0.49
1:B:314:MET:SD	1:B:329:ARG:CD	3.00	0.49
1:A:13:GLU:CD	1:A:58:ARG:HH21	2.21	0.49
1:B:44:LEU:HD13	1:B:62:TRP:HZ2	1.78	0.49
1:B:114:LEU:HD21	1:B:174:ILE:HD13	1.94	0.49
1:B:220:ALA:C	1:B:222:SER:N	2.68	0.49
1:A:336:LEU:O	1:A:336:LEU:HD23	2.13	0.48
1:B:128:VAL:HA	1:B:160:THR:O	2.13	0.48
1:B:221:SER:O	1:B:223:SER:N	2.45	0.48
1:A:26:PHE:CG	1:A:67:PHE:HB3	2.48	0.48
1:A:314:MET:SD	1:A:329:ARG:CD	3.01	0.48
1:B:25:GLN:HG3	1:B:26:PHE:N	2.28	0.48
1:A:189:LYS:O	1:A:190:GLU:C	2.56	0.48
1:B:80:ASP:O	1:B:94:GLN:HA	2.13	0.48
1:B:104:SER:CA	1:B:279:LYS:HZ3	2.24	0.48
1:B:183:SER:OG	1:B:286:GLU:HG2	2.13	0.48
1:B:190:GLU:HA	1:B:262:ALA:O	2.13	0.48
1:B:19:SER:O	1:B:21:LYS:N	2.43	0.48
1:B:37:LEU:HD12	1:B:37:LEU:C	2.39	0.48
1:B:132:SER:OG	1:B:136:LYS:HB3	2.13	0.48
1:A:303:CYS:HG	1:A:345:CYS:HG	1.55	0.48
1:B:342:MET:SD	1:B:360:LYS:HD2	2.53	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:362:LEU:HB2	1:A:363:PRO:HD3	1.96	0.48
1:A:72:LYS:O	1:A:73:ASN:C	2.56	0.48
1:B:72:LYS:O	1:B:73:ASN:C	2.56	0.48
1:B:130:CYS:CB	1:B:159:CYS:HG	2.26	0.48
1:B:162:LEU:CD2	1:B:162:LEU:C	2.87	0.48
1:B:198:PHE:HB3	1:B:285:GLN:OE1	2.14	0.48
1:B:275:ALA:HB2	1:B:282:LYS:HB3	1.96	0.48
1:B:288:ASN:HB3	1:B:308:PRO:HG3	1.94	0.48
1:A:275:ALA:HB2	1:A:282:LYS:HB3	1.96	0.48
1:A:342:MET:CG	1:A:360:LYS:HG2	2.44	0.48
1:B:59:ARG:HG2	1:B:62:TRP:CZ3	2.49	0.48
1:B:206:LYS:HB3	1:B:207:LEU:H	1.49	0.47
1:B:208:THR:O	1:B:276:LEU:HB3	2.14	0.47
1:B:333:VAL:HG13	1:B:334:TRP:N	2.29	0.47
1:A:302:THR:HG23	1:A:303:CYS:N	2.29	0.47
1:B:215:TRP:CD1	1:B:216:GLN:N	2.83	0.47
1:A:358:ASN:ND2	1:B:358:ASN:HD21	2.13	0.47
1:B:114:LEU:CD2	1:B:174:ILE:HD13	2.45	0.47
1:A:109:LEU:HD12	1:A:109:LEU:H	1.79	0.47
1:A:161:VAL:O	1:A:161:VAL:HG12	2.14	0.47
1:A:220:ALA:C	1:A:222:SER:N	2.72	0.47
1:B:100:LEU:HD12	1:B:118:LEU:CD1	2.44	0.47
1:B:290:VAL:CG1	1:B:306:TRP:O	2.62	0.47
1:A:13:GLU:OE1	1:A:58:ARG:NH2	2.48	0.47
1:A:46:LYS:HD2	1:A:52:ASN:O	2.13	0.47
1:A:102:ALA:HA	1:A:115:THR:O	2.14	0.47
1:A:177:LEU:HD21	1:A:200:LEU:HD22	1.97	0.47
1:A:183:SER:OG	1:A:286:GLU:HG2	2.15	0.47
1:B:107:HIS:ND1	1:B:278:ALA:HB3	2.28	0.47
1:B:189:LYS:HA	1:B:292:MET:H	1.79	0.47
1:B:205:GLU:HG3	1:B:208:THR:HG21	1.93	0.47
1:B:234:LYS:HA	1:B:253:LEU:CG	2.29	0.47
1:A:111:GLY:HA2	1:A:148:GLN:HB3	1.97	0.47
1:B:82:TYR:N	1:B:82:TYR:CD1	2.83	0.47
1:A:276:LEU:N	1:A:276:LEU:HD22	2.21	0.47
1:A:303:CYS:SG	1:A:345:CYS:SG	3.08	0.47
1:B:194:VAL:HG23	1:B:262:ALA:HB2	1.97	0.47
1:B:305:VAL:O	1:B:331:LYS:HB2	2.15	0.47
1:A:195:GLU:HG2	1:A:258:THR:HG22	1.97	0.46
1:A:293:ARG:O	1:A:303:CYS:HA	2.14	0.46
1:A:190:GLU:HA	1:A:262:ALA:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:LEU:O	1:A:226:TRP:HB3	2.16	0.46
1:B:130:CYS:N	1:B:138:ILE:O	2.47	0.46
1:B:115:THR:HA	1:B:145:SER:CB	2.45	0.46
1:A:141:GLY:C	1:A:143:THR:N	2.73	0.46
1:A:205:GLU:HG2	1:A:208:THR:HG21	1.98	0.46
1:B:29:LYS:HD3	1:B:33:GLN:OE1	2.16	0.46
1:A:290:VAL:HG22	1:A:308:PRO:HG2	1.97	0.46
1:A:333:VAL:HG13	1:A:334:TRP:N	2.30	0.46
1:B:98:PHE:CG	1:B:161:VAL:HG11	2.51	0.46
1:B:294:ALA:HB2	1:B:303:CYS:SG	2.56	0.46
1:A:59:ARG:HA	1:A:62:TRP:CG	2.50	0.46
1:A:59:ARG:HH11	1:A:59:ARG:HB2	1.80	0.46
1:A:86:VAL:O	1:A:88:ASP:N	2.43	0.46
1:B:189:LYS:O	1:B:190:GLU:C	2.58	0.46
1:B:194:VAL:CG2	1:B:262:ALA:HB2	2.46	0.46
1:A:29:LYS:HB3	1:A:29:LYS:HE3	1.77	0.46
1:A:202:PHE:O	1:A:203:THR:O	2.33	0.46
1:A:294:ALA:HB2	1:A:303:CYS:SG	2.56	0.46
1:A:109:LEU:HD13	1:A:281:GLY:HA3	1.97	0.46
1:A:198:PHE:HB3	1:A:285:GLN:OE1	2.15	0.46
1:B:87:GLU:O	1:B:88:ASP:CG	2.59	0.46
1:B:338:PRO:HB2	1:B:361:VAL:HG21	1.97	0.46
1:A:202:PHE:HE2	1:A:234:LYS:NZ	2.14	0.46
1:B:27:HIS:HD2	1:B:85:GLU:HB2	1.81	0.46
1:B:141:GLY:C	1:B:143:THR:H	2.22	0.46
1:A:305:VAL:HG21	1:A:347:LEU:HG	1.96	0.45
1:B:117:THR:HA	1:B:143:THR:OG1	2.15	0.45
1:A:115:THR:HA	1:A:145:SER:CB	2.46	0.45
1:A:117:THR:HA	1:A:143:THR:OG1	2.16	0.45
1:A:162:LEU:C	1:A:162:LEU:CD2	2.89	0.45
1:B:107:HIS:ND1	1:B:278:ALA:HB2	2.32	0.45
1:B:187:TYR:HB3	1:B:355:LEU:HG	1.98	0.45
1:A:129:GLN:HA	1:A:139:GLN:HB3	1.98	0.45
1:A:161:VAL:HB	1:A:168:VAL:CG1	2.46	0.45
1:A:212:GLU:HG3	1:A:214:TRP:NE1	2.30	0.45
1:A:231:LEU:C	1:A:232:LYS:HG3	2.42	0.45
1:B:36:ILE:O	1:B:47:GLY:N	2.49	0.45
1:B:110:GLN:HE21	1:B:110:GLN:HB3	1.58	0.45
1:B:196:PHE:HE2	1:B:259:LEU:HD12	1.81	0.45
1:B:219:ARG:O	1:B:221:SER:N	2.49	0.45
1:A:118:LEU:HD23	1:A:120:SER:OG	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:251:LYS:HB2	1:B:251:LYS:NZ	2.31	0.45
1:A:109:LEU:HD12	1:A:109:LEU:N	2.31	0.45
1:B:178:ALA:CB	1:B:283:LEU:HD22	2.46	0.45
1:A:14:LEU:HD22	1:A:69:LEU:HD23	1.99	0.45
1:A:46:LYS:HE2	1:A:46:LYS:HB2	1.66	0.45
1:A:344:GLN:HE22	1:B:344:GLN:CD	2.25	0.45
1:B:107:HIS:HB2	1:B:278:ALA:HB3	1.99	0.45
1:B:195:GLU:HG2	1:B:258:THR:HG22	1.99	0.45
1:A:342:MET:HG2	1:A:360:LYS:HG2	1.98	0.45
1:B:121:PRO:HA	1:B:122:PRO:HD3	1.84	0.45
1:B:213:LEU:O	1:B:226:TRP:HB3	2.16	0.45
1:A:151:LEU:O	1:A:154:SER:HB2	2.17	0.45
1:A:198:PHE:CD2	1:A:272:LEU:HD22	2.52	0.45
1:A:111:GLY:O	1:A:112:GLN:C	2.60	0.45
1:A:358:ASN:HD21	1:B:358:ASN:ND2	2.14	0.45
1:B:177:LEU:HD11	1:B:276:LEU:HD11	1.98	0.45
1:B:72:LYS:O	1:B:74:LEU:N	2.51	0.44
1:A:20:GLN:O	1:A:22:LYS:N	2.49	0.44
1:A:296:GLN:HA	1:A:301:LEU:HA	1.98	0.44
1:B:30:ASN:O	1:B:32:ASN:N	2.50	0.44
1:B:130:CYS:HB2	1:B:138:ILE:HG23	1.98	0.44
1:A:100:LEU:HD12	1:A:118:LEU:CD1	2.48	0.44
1:B:44:LEU:HD13	1:B:62:TRP:CZ2	2.52	0.44
1:A:30:ASN:HA	1:A:82:TYR:HA	2.00	0.44
1:A:267:ALA:HB1	1:A:307:GLY:CA	2.46	0.44
1:B:30:ASN:C	1:B:32:ASN:N	2.74	0.44
1:B:52:ASN:OD1	1:B:53:ASP:N	2.50	0.44
1:B:77:GLU:H	1:B:77:GLU:HG2	1.63	0.44
1:B:315:LEU:HD22	1:B:316:SER:H	1.82	0.44
1:A:37:LEU:C	1:A:37:LEU:HD12	2.43	0.44
1:A:259:LEU:HD13	1:A:289:LEU:HD22	1.99	0.44
1:B:116:LEU:O	1:B:143:THR:HA	2.18	0.44
1:B:298:GLN:HB2	1:B:299:LYS:HD3	2.00	0.44
1:B:8:LYS:C	1:B:10:ASP:H	2.26	0.44
1:B:326:VAL:HG11	1:B:335:VAL:HG22	2.00	0.44
1:A:114:LEU:HD21	1:A:174:ILE:HD13	1.99	0.44
1:A:214:TRP:CE3	1:A:226:TRP:HZ3	2.34	0.44
1:B:28:TRP:CE2	1:B:84:CYS:SG	3.03	0.44
1:A:26:PHE:CD1	1:A:67:PHE:CD1	3.06	0.44
1:A:82:TYR:N	1:A:82:TYR:CD1	2.86	0.44
1:A:128:VAL:HA	1:A:160:THR:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:303:CYS:HB3	1:B:345:CYS:HG	1.83	0.44
1:A:28:TRP:CE2	1:A:84:CYS:SG	2.98	0.43
1:A:187:TYR:HB3	1:A:355:LEU:HG	2.00	0.43
1:B:89:GLN:HE21	1:B:89:GLN:HB3	1.65	0.43
1:B:309:THR:HB	1:B:313:LEU:CD1	2.40	0.43
1:B:333:VAL:CG1	1:B:334:TRP:N	2.81	0.43
1:B:36:ILE:HD12	1:B:51:LEU:HD23	2.01	0.43
1:B:263:LEU:O	1:B:266:TYR:HD1	2.00	0.43
1:A:109:LEU:HB3	1:A:283:LEU:CD2	2.48	0.43
1:A:293:ARG:C	1:A:303:CYS:SG	3.02	0.43
1:B:30:ASN:C	1:B:32:ASN:H	2.27	0.43
1:B:130:CYS:O	1:B:138:ILE:HG22	2.18	0.43
1:B:177:LEU:CD2	1:B:283:LEU:HD11	2.29	0.43
1:A:215:TRP:CD1	1:A:216:GLN:N	2.87	0.43
1:B:162:LEU:CD2	1:B:163:GLN:N	2.78	0.43
1:B:198:PHE:CD2	1:B:272:LEU:HD22	2.54	0.43
1:B:231:LEU:C	1:B:232:LYS:HG3	2.43	0.43
1:A:44:LEU:HD13	1:A:62:TRP:CZ2	2.53	0.43
1:A:121:PRO:HA	1:A:122:PRO:HD3	1.93	0.43
1:A:128:VAL:HB	1:A:144:LEU:HD23	2.00	0.43
1:B:17:THR:HG23	1:B:65:GLY:C	2.44	0.43
1:B:27:HIS:HB2	1:B:35:LYS:HZ3	1.84	0.43
1:B:51:LEU:HD12	1:B:51:LEU:HA	1.77	0.43
1:A:98:PHE:CG	1:A:161:VAL:HG11	2.54	0.43
1:B:13:GLU:CD	1:B:58:ARG:HH21	2.25	0.43
1:B:26:PHE:CD2	1:B:67:PHE:HB3	2.54	0.43
1:B:319:LEU:HA	1:B:343:TRP:HA	2.00	0.43
1:B:212:GLU:HG3	1:B:214:TRP:NE1	2.34	0.43
1:A:59:ARG:HG2	1:A:62:TRP:CE3	2.54	0.43
1:A:212:GLU:HB2	1:A:227:ILE:O	2.19	0.43
1:B:17:THR:HG22	1:B:18:ALA:N	2.33	0.43
1:A:206:LYS:HB3	1:A:207:LEU:HG	2.01	0.43
1:B:164:ASN:C	1:B:166:LYS:H	2.25	0.43
1:A:206:LYS:O	1:A:207:LEU:C	2.62	0.42
1:A:277:GLU:CG	1:A:278:ALA:N	2.59	0.42
1:A:288:ASN:HB3	1:A:308:PRO:HG3	2.01	0.42
1:B:189:LYS:O	1:B:191:GLY:N	2.52	0.42
1:B:219:ARG:NH1	1:B:306:TRP:HZ3	2.16	0.42
1:B:296:GLN:HA	1:B:301:LEU:HA	2.01	0.42
1:A:226:TRP:HD1	1:A:241:VAL:HG11	1.84	0.42
1:B:161:VAL:HB	1:B:168:VAL:CG1	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:TRP:CD1	1:A:214:TRP:N	2.87	0.42
1:B:109:LEU:CD1	1:B:280:THR:C	2.92	0.42
1:B:130:CYS:SG	1:B:144:LEU:HD21	2.59	0.42
1:A:114:LEU:HB2	1:A:149:LEU:HD11	2.01	0.42
1:A:317:LEU:HB2	1:A:326:VAL:HG12	2.02	0.42
1:B:297:LEU:N	1:B:300:ASN:O	2.53	0.42
1:B:29:LYS:HB3	1:B:29:LYS:HE3	1.97	0.42
1:B:111:GLY:HA2	1:B:148:GLN:CB	2.49	0.42
1:B:296:GLN:HG3	1:B:301:LEU:HB3	2.02	0.42
1:B:315:LEU:HD23	1:B:346:LEU:O	2.18	0.42
1:A:128:VAL:HB	1:A:144:LEU:CD2	2.50	0.42
1:A:342:MET:SD	1:A:360:LYS:CG	3.08	0.42
1:B:112:GLN:O	1:B:149:LEU:HD12	2.19	0.42
1:B:162:LEU:HD23	1:B:162:LEU:C	2.45	0.42
1:B:207:LEU:HG	1:B:207:LEU:H	1.41	0.42
1:A:253:LEU:HD22	1:A:253:LEU:HA	1.86	0.42
1:B:109:LEU:HD12	1:B:109:LEU:H	1.85	0.42
1:A:344:GLN:NE2	1:B:344:GLN:CD	2.77	0.42
1:B:104:SER:O	1:B:279:LYS:HD2	2.20	0.42
1:A:20:GLN:HG3	1:A:24:ILE:HD13	2.02	0.42
1:B:214:TRP:CE3	1:B:226:TRP:HZ3	2.34	0.42
1:B:362:LEU:CB	1:B:363:PRO:HD3	2.49	0.42
1:A:110:GLN:NE2	1:A:178:ALA:HB2	2.33	0.42
1:A:294:ALA:N	1:A:303:CYS:SG	2.93	0.42
1:B:54:ARG:HG2	1:B:72:LYS:HB2	2.01	0.42
1:B:133:PRO:HG3	1:B:156:THR:O	2.19	0.42
1:A:51:LEU:HD12	1:A:51:LEU:HA	1.73	0.41
1:B:247:LEU:H	1:B:247:LEU:HD23	1.85	0.41
1:A:130:CYS:SG	1:A:144:LEU:HD11	2.61	0.41
1:B:30:ASN:C	1:B:30:ASN:OD1	2.62	0.41
1:A:267:ALA:HB1	1:A:307:GLY:HA3	2.01	0.41
1:B:20:GLN:O	1:B:22:LYS:N	2.54	0.41
1:A:52:ASN:CG	1:A:53:ASP:N	2.77	0.41
1:A:296:GLN:HA	1:A:301:LEU:CB	2.51	0.41
1:B:196:PHE:HD2	1:B:257:LEU:O	2.03	0.41
1:A:14:LEU:CD2	1:A:69:LEU:HD23	2.50	0.41
1:A:194:VAL:HG23	1:A:262:ALA:HB2	2.02	0.41
1:A:315:LEU:HD23	1:A:346:LEU:O	2.20	0.41
1:A:344:GLN:CD	1:B:344:GLN:NE2	2.78	0.41
1:B:7:LYS:HB2	1:B:10:ASP:CB	2.50	0.41
1:B:16:CYS:SG	1:B:67:PHE:HB2	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:299:LYS:HD3	1:B:299:LYS:N	2.29	0.41
1:A:100:LEU:HA	1:A:118:LEU:HD12	2.03	0.41
1:B:46:LYS:HB2	1:B:46:LYS:HE2	1.72	0.41
1:B:349:ASP:O	1:B:351:GLY:N	2.51	0.41
1:A:42:SER:HG	1:A:43:PHE:HD1	1.66	0.41
1:A:263:LEU:HB2	1:A:266:TYR:CE1	2.56	0.41
1:A:302:THR:CG2	1:A:303:CYS:N	2.84	0.41
1:B:87:GLU:O	1:B:88:ASP:CB	2.69	0.41
1:B:128:VAL:HB	1:B:144:LEU:CD2	2.51	0.41
1:B:293:ARG:O	1:B:303:CYS:HA	2.20	0.41
1:A:100:LEU:HD22	1:A:159:CYS:HB2	2.03	0.41
1:B:302:THR:HG23	1:B:303:CYS:N	2.35	0.41
1:A:175:VAL:O	1:A:175:VAL:HG12	2.20	0.41
1:A:186:VAL:HG12	1:A:188:LYS:HG2	2.02	0.41
1:A:208:THR:O	1:A:276:LEU:HB2	2.19	0.41
1:A:300:ASN:HA	1:A:338:PRO:HD3	2.03	0.41
1:A:349:ASP:O	1:A:351:GLY:N	2.53	0.41
1:B:136:LYS:HD2	1:B:138:ILE:HB	2.02	0.41
1:B:180:GLN:O	1:B:180:GLN:HG3	2.20	0.41
1:A:58:ARG:NE	1:A:61:LEU:HD12	2.36	0.41
1:B:51:LEU:O	1:B:52:ASN:C	2.64	0.41
1:B:86:VAL:HG12	1:B:87:GLU:H	1.86	0.41
1:B:114:LEU:HB2	1:B:149:LEU:HD11	2.03	0.41
1:B:205:GLU:CG	1:B:208:THR:CG2	2.93	0.41
1:B:300:ASN:OD1	1:B:337:ASN:N	2.54	0.40
1:B:2:LYS:H	1:B:2:LYS:HG3	1.72	0.40
1:B:246:LYS:HG3	1:B:246:LYS:O	2.21	0.40
1:B:291:VAL:HG12	1:B:292:MET:N	2.36	0.40
1:B:343:TRP:N	1:B:359:ILE:O	2.53	0.40
1:A:315:LEU:CD1	1:A:333:VAL:HB	2.51	0.40
1:A:319:LEU:HB3	1:A:343:TRP:NE1	2.36	0.40
1:B:241:VAL:O	1:B:242:THR:OG1	2.34	0.40
1:A:51:LEU:O	1:A:52:ASN:C	2.65	0.40
1:A:347:LEU:O	1:A:354:LEU:HB2	2.20	0.40
1:B:202:PHE:HE2	1:B:234:LYS:NZ	2.20	0.40
1:B:208:THR:O	1:B:276:LEU:HB2	2.21	0.40
1:B:215:TRP:HD1	1:B:216:GLN:C	2.30	0.40
1:A:356:GLU:OE2	1:B:344:GLN:OE1	2.39	0.40
1:B:59:ARG:HA	1:B:62:TRP:CD2	2.57	0.40
1:B:100:LEU:HD12	1:B:118:LEU:HD12	2.02	0.40
1:B:296:GLN:HA	1:B:301:LEU:CB	2.52	0.40

All (19) 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:9:GLY:O	1:B:73:ASN:OD1[2_867]	0.59	1.61
1:A:73:ASN:ND2	1:B:73:ASN:CA[2_867]	0.99	1.21
1:A:73:ASN:OD1	1:B:9:GLY:O[2_867]	1.06	1.14
1:A:73:ASN:OD1	1:B:9:GLY:C[2_867]	1.27	0.93
1:A:9:GLY:C	1:B:73:ASN:OD1[2_867]	1.58	0.62
1:A:73:ASN:CA	1:B:73:ASN:ND2[2_867]	1.59	0.61
1:A:73:ASN:OD1	1:B:9:GLY:CA[2_867]	1.61	0.59
1:A:9:GLY:O	1:B:73:ASN:CG[2_867]	1.70	0.50
1:A:204:VAL:CG2	1:B:25:GLN:OE1[2_867]	1.72	0.48
1:A:44:LEU:O	1:B:173:ASP:OD2[2_867]	1.76	0.44
1:A:73:ASN:N	1:B:73:ASN:ND2[2_867]	1.78	0.42
1:A:73:ASN:ND2	1:B:73:ASN:N[2_867]	1.84	0.36
1:A:25:GLN:OE1	1:B:204:VAL:N[2_867]	1.86	0.34
1:A:73:ASN:ND2	1:B:73:ASN:CB[2_867]	1.97	0.23
1:A:73:ASN:CG	1:B:9:GLY:O[2_867]	1.98	0.22
1:A:106:THR:OG1	1:B:48:PRO:CD[2_867]	2.08	0.12
1:A:35:LYS:CE	1:B:205:GLU:OE1[2_867]	2.10	0.10
1:A:73:ASN:CG	1:B:73:ASN:CA[2_867]	2.13	0.07
1:A:48:PRO:CA	1:B:105:ASP:OD1[2_867]	2.19	0.01

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	361/363 (99%)	259 (72%)	58 (16%)	44 (12%)	0	4
1	B	361/363 (99%)	260 (72%)	59 (16%)	42 (12%)	0	5
All	All	722/726 (99%)	519 (72%)	117 (16%)	86 (12%)	0	5

All (86) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	21	LYS
1	A	52	ASN
1	A	73	ASN
1	A	87	GLU
1	A	88	ASP
1	A	105	ASP
1	A	136	LYS
1	A	147	SER
1	A	151	LEU
1	A	178	ALA
1	A	179	PHE
1	A	190	GLU
1	A	202	PHE
1	A	203	THR
1	A	205	GLU
1	A	207	LEU
1	A	222	SER
1	A	241	VAL
1	A	242	THR
1	A	244	ASP
1	A	255	LEU
1	A	260	PRO
1	A	280	THR
1	A	349	ASP
1	B	18	ALA
1	B	21	LYS
1	B	52	ASN
1	B	73	ASN
1	B	87	GLU
1	B	88	ASP
1	B	105	ASP
1	B	147	SER
1	B	151	LEU
1	B	190	GLU
1	B	202	PHE
1	B	205	GLU
1	B	241	VAL
1	B	242	THR
1	B	244	ASP
1	B	255	LEU
1	B	261	GLN
1	B	280	THR
1	B	349	ASP

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Mol	Chain	Res	Type
1	A	53	ASP
1	A	106	THR
1	A	112	GLN
1	A	278	ALA
1	B	10	ASP
1	B	106	THR
1	B	136	LYS
1	B	178	ALA
1	B	206	LYS
1	B	207	LEU
1	B	222	SER
1	B	243	GLN
1	B	278	ALA
1	A	10	ASP
1	A	19	SER
1	A	20	GLN
1	A	233	ASN
1	A	253	LEU
1	A	297	LEU
1	A	324	ALA
1	B	20	GLN
1	B	53	ASP
1	B	112	GLN
1	B	220	ALA
1	B	240	ARG
1	B	253	LEU
1	B	324	ALA
1	B	351	GLY
1	A	16	CYS
1	A	201	ALA
1	A	206	LYS
1	A	240	ARG
1	A	351	GLY
1	B	16	CYS
1	B	23	SER
1	B	201	ALA
1	B	326	VAL
1	A	113	SER
1	A	220	ALA
1	A	326	VAL
1	B	126	PRO
1	A	362	LEU

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Mol	Chain	Res	Type
1	B	122	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	318/324 (98%)	216 (68%)	102 (32%)	0	2
1	B	318/324 (98%)	223 (70%)	95 (30%)	0	3
All	All	636/648 (98%)	439 (69%)	197 (31%)	0	2

All (197) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	3	VAL
1	A	11	THR
1	A	14	LEU
1	A	15	THR
1	A	16	CYS
1	A	19	SER
1	A	23	SER
1	A	25	GLN
1	A	29	LYS
1	A	35	LYS
1	A	36	ILE
1	A	43	PHE
1	A	46	LYS
1	A	59	ARG
1	A	63	ASP
1	A	68	PRO
1	A	70	ILE
1	A	76	ILE
1	A	84	CYS
1	A	89	GLN
1	A	97	VAL

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Mol	Chain	Res	Type
1	A	101	THR
1	A	103	ASN
1	A	104	SER
1	A	115	THR
1	A	118	LEU
1	A	124	SER
1	A	125	SER
1	A	132	SER
1	A	134	ARG
1	A	136	LYS
1	A	139	GLN
1	A	143	THR
1	A	144	LEU
1	A	148	GLN
1	A	151	LEU
1	A	152	GLN
1	A	154	SER
1	A	156	THR
1	A	158	THR
1	A	163	GLN
1	A	164	ASN
1	A	166	LYS
1	A	171	LYS
1	A	175	VAL
1	A	185	ILE
1	A	193	GLN
1	A	195	GLU
1	A	197	SER
1	A	198	PHE
1	A	207	LEU
1	A	208	THR
1	A	216	GLN
1	A	221	SER
1	A	223	SER
1	A	225	SER
1	A	226	TRP
1	A	228	THR
1	A	229	PHE
1	A	230	ASP
1	A	234	LYS
1	A	235	GLU
1	A	237	SER

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Mol	Chain	Res	Type
1	A	238	VAL
1	A	243	GLN
1	A	247	LEU
1	A	251	LYS
1	A	253	LEU
1	A	257	LEU
1	A	259	LEU
1	A	260	PRO
1	A	266	TYR
1	A	272	LEU
1	A	273	THR
1	A	274	LEU
1	A	276	LEU
1	A	279	LYS
1	A	282	LYS
1	A	285	GLN
1	A	286	GLU
1	A	295	THR
1	A	298	GLN
1	A	299	LYS
1	A	302	THR
1	A	304	GLU
1	A	305	VAL
1	A	310	SER
1	A	312	LYS
1	A	313	LEU
1	A	315	LEU
1	A	316	SER
1	A	320	GLU
1	A	329	ARG
1	A	330	GLU
1	A	333	VAL
1	A	336	LEU
1	A	339	GLU
1	A	342	MET
1	A	350	SER
1	A	353	VAL
1	A	362	LEU
1	B	3	VAL
1	B	11	THR
1	B	12	VAL
1	B	14	LEU

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Mol	Chain	Res	Type
1	B	15	THR
1	B	19	SER
1	B	23	SER
1	B	25	GLN
1	B	29	LYS
1	B	35	LYS
1	B	43	PHE
1	B	58	ARG
1	B	59	ARG
1	B	63	ASP
1	B	68	PRO
1	B	70	ILE
1	B	76	ILE
1	B	82	TYR
1	B	89	GLN
1	B	97	VAL
1	B	101	THR
1	B	103	ASN
1	B	115	THR
1	B	118	LEU
1	B	125	SER
1	B	127	SER
1	B	132	SER
1	B	134	ARG
1	B	139	GLN
1	B	143	THR
1	B	144	LEU
1	B	148	GLN
1	B	151	LEU
1	B	152	GLN
1	B	154	SER
1	B	156	THR
1	B	158	THR
1	B	162	LEU
1	B	163	GLN
1	B	164	ASN
1	B	166	LYS
1	B	175	VAL
1	B	185	ILE
1	B	193	GLN
1	B	195	GLU
1	B	197	SER

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Mol	Chain	Res	Type
1	B	198	PHE
1	B	203	THR
1	B	207	LEU
1	B	208	THR
1	B	216	GLN
1	B	221	SER
1	B	223	SER
1	B	225	SER
1	B	226	TRP
1	B	228	THR
1	B	229	PHE
1	B	230	ASP
1	B	235	GLU
1	B	236	VAL
1	B	237	SER
1	B	238	VAL
1	B	243	GLN
1	B	247	LEU
1	B	251	LYS
1	B	253	LEU
1	B	257	LEU
1	B	259	LEU
1	B	273	THR
1	B	274	LEU
1	B	276	LEU
1	B	279	LYS
1	B	282	LYS
1	B	285	GLN
1	B	286	GLU
1	B	298	GLN
1	B	299	LYS
1	B	302	THR
1	B	304	GLU
1	B	305	VAL
1	B	310	SER
1	B	312	LYS
1	B	313	LEU
1	B	315	LEU
1	B	316	SER
1	B	320	GLU
1	B	329	ARG
1	B	330	GLU

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Mol	Chain	Res	Type
1	B	333	VAL
1	B	336	LEU
1	B	339	GLU
1	B	342	MET
1	B	350	SER
1	B	353	VAL
1	B	362	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	GLN
1	A	66	ASN
1	A	89	GLN
1	A	103	ASN
1	A	110	GLN
1	A	163	GLN
1	A	180	GLN
1	A	193	GLN
1	A	261	GLN
1	A	284	HIS
1	A	288	ASN
1	A	321	ASN
1	A	344	GLN
1	A	358	ASN
1	B	20	GLN
1	B	66	ASN
1	B	89	GLN
1	B	103	ASN
1	B	110	GLN
1	B	163	GLN
1	B	261	GLN
1	B	284	HIS
1	B	288	ASN
1	B	344	GLN
1	B	358	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.