



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:28 AM UTC

PDB ID : 3BQ4 / pdb_00003bq4
Title : Crystal Structure of Ad35 fiber knob
Authors : Pache, L.; Venkataraman, S.; Nemerow, G.R.; Reddy, V.S.
Deposited on : 2007-12-19
Resolution : 2.70 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:

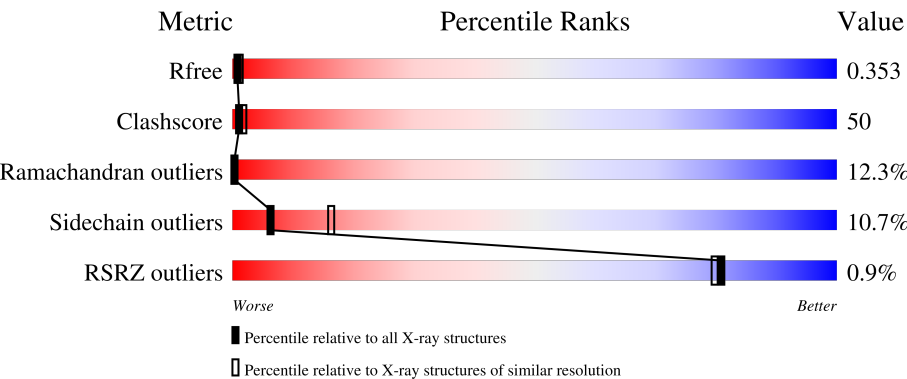
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	220	% 32%36%17%•13%
1	B	220	26%43%16%•13%
1	D	220	% 27%43%15%•13%
1	E	220	% 30%40%16%•13%
1	F	220	% 29%39%16%•13%

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Mol	Chain	Length	Quality of chain
1	G	220	30% 42% 15% 13%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 8976 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 Fiber.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	192	Total	C	N	O	S	0	0	0
			1496	944	240	305	7			
1	B	192	Total	C	N	O	S	0	0	0
			1496	944	240	305	7			
1	D	192	Total	C	N	O	S	0	0	0
			1496	944	240	305	7			
1	E	192	Total	C	N	O	S	0	0	0
			1496	944	240	305	7			
1	F	192	Total	C	N	O	S	0	0	0
			1496	944	240	305	7			
1	G	192	Total	C	N	O	S	0	0	0
			1496	944	240	305	7			

There are 102 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	104	GLY	-	expression tag	UNP Q7T925
A	105	SER	-	expression tag	UNP Q7T925
A	106	HIS	-	expression tag	UNP Q7T925
A	107	MET	-	expression tag	UNP Q7T925
A	108	ALA	-	expression tag	UNP Q7T925
A	109	SER	-	expression tag	UNP Q7T925
A	110	MET	-	expression tag	UNP Q7T925
A	111	THR	-	expression tag	UNP Q7T925
A	112	GLY	-	expression tag	UNP Q7T925
A	113	GLY	-	expression tag	UNP Q7T925
A	114	GLN	-	expression tag	UNP Q7T925
A	115	GLN	-	expression tag	UNP Q7T925
A	116	MET	-	expression tag	UNP Q7T925
A	117	GLY	-	expression tag	UNP Q7T925
A	118	ARG	-	expression tag	UNP Q7T925
A	119	GLY	-	expression tag	UNP Q7T925
A	120	SER	-	expression tag	UNP Q7T925

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Chain	Residue	Modelled	Actual	Comment	Reference
B	104	GLY	-	expression tag	UNP Q7T925
B	105	SER	-	expression tag	UNP Q7T925
B	106	HIS	-	expression tag	UNP Q7T925
B	107	MET	-	expression tag	UNP Q7T925
B	108	ALA	-	expression tag	UNP Q7T925
B	109	SER	-	expression tag	UNP Q7T925
B	110	MET	-	expression tag	UNP Q7T925
B	111	THR	-	expression tag	UNP Q7T925
B	112	GLY	-	expression tag	UNP Q7T925
B	113	GLY	-	expression tag	UNP Q7T925
B	114	GLN	-	expression tag	UNP Q7T925
B	115	GLN	-	expression tag	UNP Q7T925
B	116	MET	-	expression tag	UNP Q7T925
B	117	GLY	-	expression tag	UNP Q7T925
B	118	ARG	-	expression tag	UNP Q7T925
B	119	GLY	-	expression tag	UNP Q7T925
B	120	SER	-	expression tag	UNP Q7T925
D	104	GLY	-	expression tag	UNP Q7T925
D	105	SER	-	expression tag	UNP Q7T925
D	106	HIS	-	expression tag	UNP Q7T925
D	107	MET	-	expression tag	UNP Q7T925
D	108	ALA	-	expression tag	UNP Q7T925
D	109	SER	-	expression tag	UNP Q7T925
D	110	MET	-	expression tag	UNP Q7T925
D	111	THR	-	expression tag	UNP Q7T925
D	112	GLY	-	expression tag	UNP Q7T925
D	113	GLY	-	expression tag	UNP Q7T925
D	114	GLN	-	expression tag	UNP Q7T925
D	115	GLN	-	expression tag	UNP Q7T925
D	116	MET	-	expression tag	UNP Q7T925
D	117	GLY	-	expression tag	UNP Q7T925
D	118	ARG	-	expression tag	UNP Q7T925
D	119	GLY	-	expression tag	UNP Q7T925
D	120	SER	-	expression tag	UNP Q7T925
E	104	GLY	-	expression tag	UNP Q7T925
E	105	SER	-	expression tag	UNP Q7T925
E	106	HIS	-	expression tag	UNP Q7T925
E	107	MET	-	expression tag	UNP Q7T925
E	108	ALA	-	expression tag	UNP Q7T925
E	109	SER	-	expression tag	UNP Q7T925
E	110	MET	-	expression tag	UNP Q7T925
E	111	THR	-	expression tag	UNP Q7T925

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Chain	Residue	Modelled	Actual	Comment	Reference
E	112	GLY	-	expression tag	UNP Q7T925
E	113	GLY	-	expression tag	UNP Q7T925
E	114	GLN	-	expression tag	UNP Q7T925
E	115	GLN	-	expression tag	UNP Q7T925
E	116	MET	-	expression tag	UNP Q7T925
E	117	GLY	-	expression tag	UNP Q7T925
E	118	ARG	-	expression tag	UNP Q7T925
E	119	GLY	-	expression tag	UNP Q7T925
E	120	SER	-	expression tag	UNP Q7T925
F	104	GLY	-	expression tag	UNP Q7T925
F	105	SER	-	expression tag	UNP Q7T925
F	106	HIS	-	expression tag	UNP Q7T925
F	107	MET	-	expression tag	UNP Q7T925
F	108	ALA	-	expression tag	UNP Q7T925
F	109	SER	-	expression tag	UNP Q7T925
F	110	MET	-	expression tag	UNP Q7T925
F	111	THR	-	expression tag	UNP Q7T925
F	112	GLY	-	expression tag	UNP Q7T925
F	113	GLY	-	expression tag	UNP Q7T925
F	114	GLN	-	expression tag	UNP Q7T925
F	115	GLN	-	expression tag	UNP Q7T925
F	116	MET	-	expression tag	UNP Q7T925
F	117	GLY	-	expression tag	UNP Q7T925
F	118	ARG	-	expression tag	UNP Q7T925
F	119	GLY	-	expression tag	UNP Q7T925
F	120	SER	-	expression tag	UNP Q7T925
G	104	GLY	-	expression tag	UNP Q7T925
G	105	SER	-	expression tag	UNP Q7T925
G	106	HIS	-	expression tag	UNP Q7T925
G	107	MET	-	expression tag	UNP Q7T925
G	108	ALA	-	expression tag	UNP Q7T925
G	109	SER	-	expression tag	UNP Q7T925
G	110	MET	-	expression tag	UNP Q7T925
G	111	THR	-	expression tag	UNP Q7T925
G	112	GLY	-	expression tag	UNP Q7T925
G	113	GLY	-	expression tag	UNP Q7T925
G	114	GLN	-	expression tag	UNP Q7T925
G	115	GLN	-	expression tag	UNP Q7T925
G	116	MET	-	expression tag	UNP Q7T925
G	117	GLY	-	expression tag	UNP Q7T925
G	118	ARG	-	expression tag	UNP Q7T925
G	119	GLY	-	expression tag	UNP Q7T925

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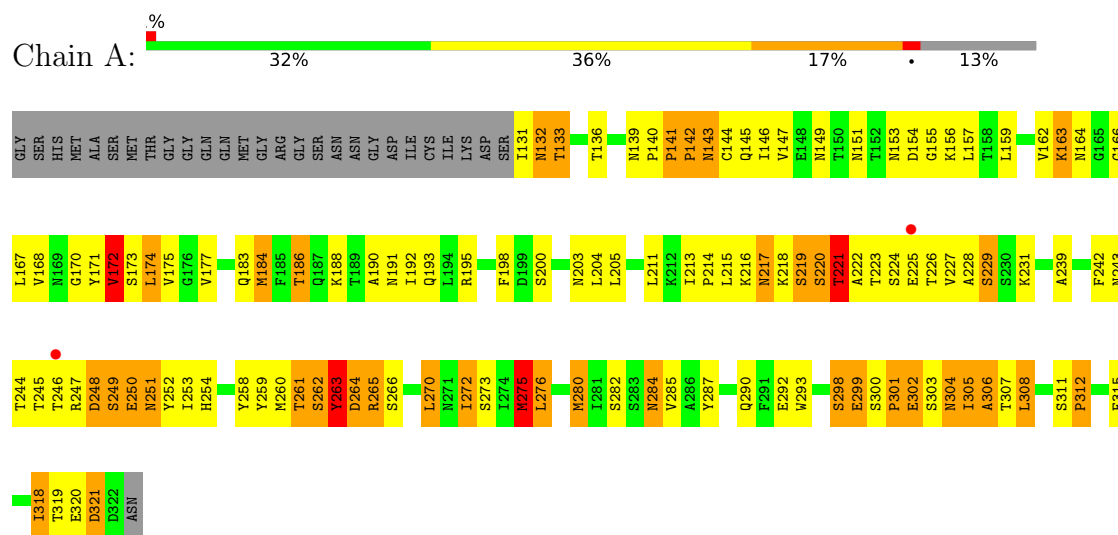
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Chain	Residue	Modelled	Actual	Comment	Reference
G	120	SER	-	expression tag	UNP Q7T925

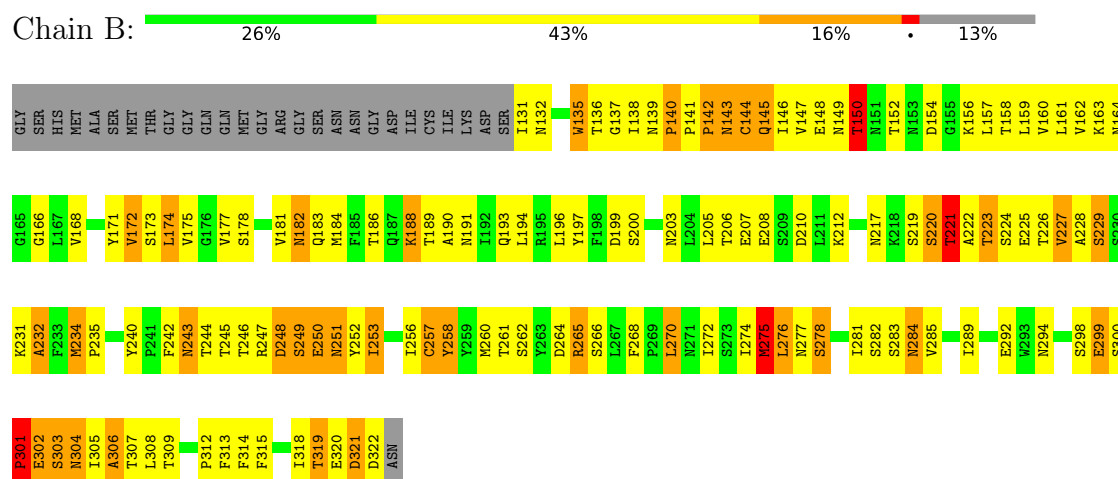
3 Residue-property plots

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: Fiber

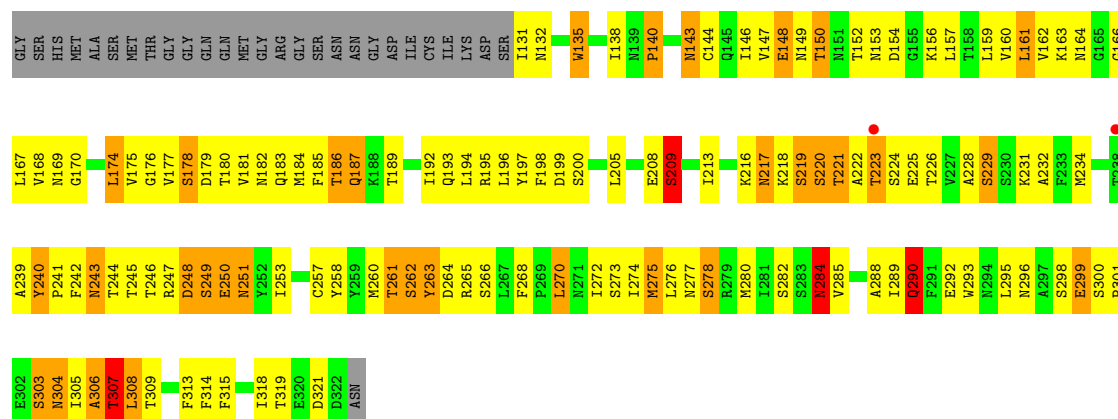


• Molecule 1: Fiber

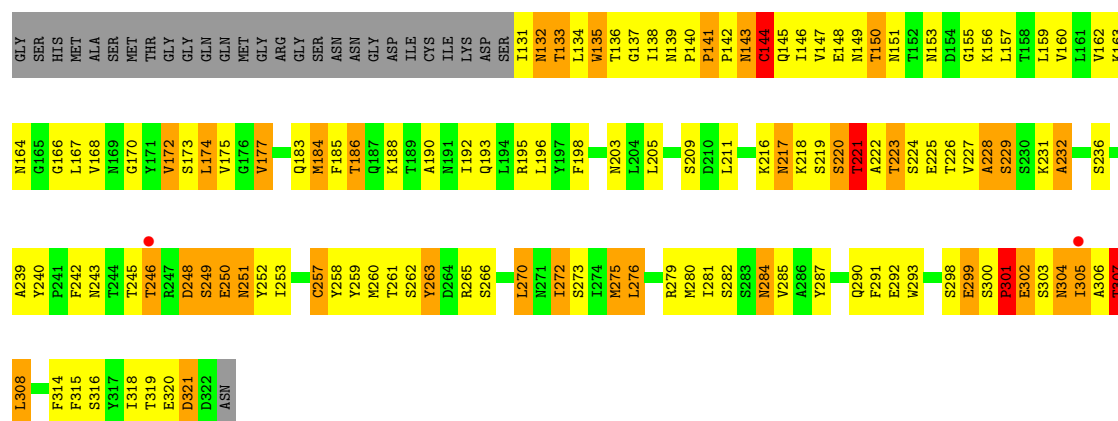


• Molecule 1: Fiber

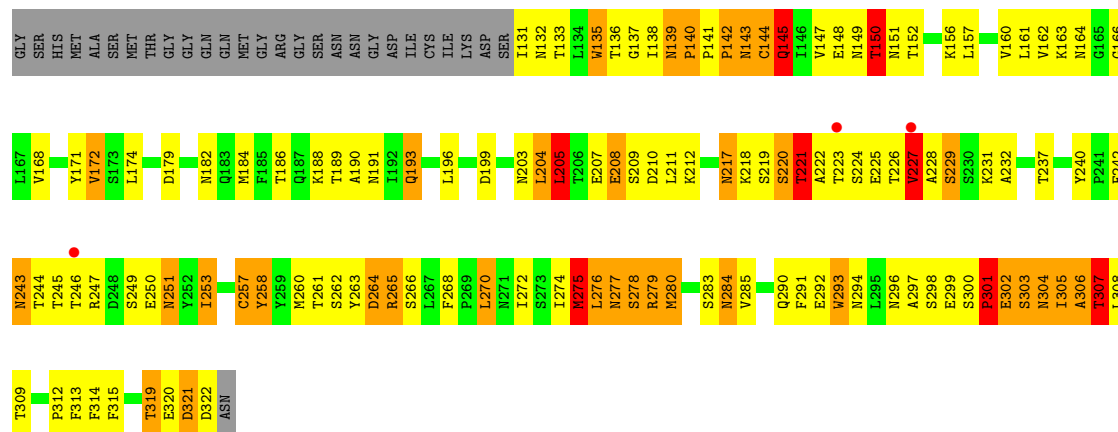




• Molecule 1: Fiber



• Molecule 1: Fiber



• Molecule 1: Fiber





4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	173.91Å 173.91Å 154.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.03 – 2.70 46.03 – 2.70	Depositor EDS
% Data completeness (in resolution range)	91.5 (46.03-2.70) 91.5 (46.03-2.70)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.69Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.332 , 0.352 0.332 , 0.353	Depositor DCC
R_{free} test set	2911 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	62.8	Xtriage
Anisotropy	0.829	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 90.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.468 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	8976	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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	0.64	1/1527 (0.1%)	1.15	11/2084 (0.5%)
1	B	0.61	0/1527	1.19	17/2084 (0.8%)
1	D	0.60	1/1527 (0.1%)	1.09	8/2084 (0.4%)
1	E	0.64	1/1527 (0.1%)	1.15	13/2084 (0.6%)
1	F	0.61	0/1527	1.25	24/2084 (1.2%)
1	G	0.60	2/1527 (0.1%)	1.07	5/2084 (0.2%)
All	All	0.62	5/9162 (0.1%)	1.15	78/12504 (0.6%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	184	MET	SD-CE	6.46	1.95	1.79
1	G	223	THR	CA-CB	5.99	1.62	1.53
1	G	175	VAL	CA-CB	5.58	1.59	1.53
1	A	184	MET	SD-CE	5.04	1.92	1.79
1	D	223	THR	CA-CB	5.02	1.61	1.53

All (78) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	221	THR	N-CA-C	-9.47	90.62	110.80
1	F	275	MET	N-CA-C	9.13	123.98	113.02
1	A	221	THR	N-CA-C	-8.99	91.64	110.80
1	F	307	THR	N-CA-C	-8.92	102.58	113.19
1	A	186	THR	N-CA-C	-8.40	103.04	113.28
1	G	221	THR	N-CA-C	-7.89	94.00	110.80
1	E	307	THR	N-CA-C	-7.78	103.80	113.28
1	D	221	THR	N-CA-C	-7.74	94.31	110.80
1	E	151	ASN	N-CA-C	-7.58	103.59	112.92
1	E	177	VAL	CB-CA-C	-7.53	104.35	111.05
1	F	144	CYS	N-CA-C	7.50	119.96	109.31
1	G	135	TRP	N-CA-C	7.46	118.37	108.07
1	A	266	SER	N-CA-C	7.38	120.38	110.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	275	MET	N-CA-C	7.34	121.99	112.89
1	B	221	THR	N-CA-C	-7.06	95.76	110.80
1	F	221	THR	N-CA-C	-7.05	95.78	110.80
1	D	135	TRP	N-CA-C	7.04	117.28	108.19
1	A	151	ASN	N-CA-C	-6.84	103.52	112.41
1	B	240	TYR	CA-C-N	6.69	126.63	120.21
1	B	240	TYR	C-N-CA	6.69	126.63	120.21
1	D	219	SER	N-CA-C	-6.68	103.21	112.30
1	B	258	TYR	N-CA-C	6.63	119.25	109.24
1	F	151	ASN	N-CA-C	-6.60	104.17	111.36
1	E	144	CYS	N-CA-C	6.44	119.20	109.41
1	F	240	TYR	CA-C-N	6.38	126.74	119.92
1	F	240	TYR	C-N-CA	6.38	126.74	119.92
1	E	266	SER	N-CA-C	6.30	119.60	110.28
1	A	133	THR	N-CA-C	6.21	118.53	108.34
1	B	210	ASP	N-CA-C	-6.18	105.38	113.17
1	B	257	CYS	N-CA-C	-6.13	100.48	109.18
1	B	136	THR	N-CA-C	-6.11	105.99	113.50
1	F	135	TRP	N-CA-C	6.10	116.61	108.38
1	F	193	GLN	N-CA-C	6.08	118.32	108.34
1	A	311	SER	N-CA-C	-6.03	101.43	110.24
1	G	238	THR	N-CA-C	-6.02	106.10	113.50
1	F	257	CYS	N-CA-C	-6.01	100.27	109.41
1	B	253	ILE	N-CA-C	-5.99	99.49	108.12
1	F	264	ASP	N-CA-C	-5.99	102.46	110.55
1	E	133	THR	N-CA-C	5.97	118.13	107.80
1	F	279	ARG	N-CA-C	-5.97	100.07	108.96
1	F	212	LYS	N-CA-C	5.96	120.68	113.17
1	B	135	TRP	N-CA-C	5.89	116.02	108.24
1	B	144	CYS	N-CA-C	5.88	118.39	109.63
1	F	208	GLU	N-CA-C	5.78	119.59	112.54
1	F	262	SER	N-CA-C	5.77	117.24	111.07
1	B	234	MET	CA-C-N	5.72	127.00	119.84
1	B	234	MET	C-N-CA	5.72	127.00	119.84
1	E	141	PRO	N-CA-C	-5.71	103.73	110.70
1	F	277	ASN	N-CA-C	5.69	119.10	111.24
1	G	239	ALA	N-CA-C	-5.66	106.04	113.12
1	A	318	ILE	CB-CA-C	-5.65	105.12	111.23
1	A	172	VAL	CB-CA-C	-5.65	102.13	110.82
1	F	258	TYR	N-CA-C	5.61	117.94	109.41
1	E	240	TYR	CA-C-N	5.58	125.57	120.21
1	E	240	TYR	C-N-CA	5.58	125.57	120.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	297	ALA	N-CA-C	5.57	113.34	108.78
1	D	240	TYR	CA-C-N	5.54	125.85	119.92
1	D	240	TYR	C-N-CA	5.54	125.85	119.92
1	E	135	TRP	N-CA-C	5.52	117.37	109.14
1	A	141	PRO	N-CA-C	-5.49	104.00	110.70
1	E	232	ALA	N-CA-C	-5.47	106.56	113.18
1	B	208	GLU	N-CA-C	5.37	119.09	112.54
1	B	319	THR	N-CA-C	5.36	117.64	110.35
1	D	290	GLN	N-CA-C	5.34	117.61	108.90
1	F	293	TRP	N-CA-C	-5.32	101.04	109.76
1	B	234	MET	N-CA-C	5.26	116.51	109.83
1	E	186	THR	N-CA-C	-5.26	106.82	113.18
1	F	210	ASP	N-CA-C	-5.17	106.97	113.28
1	F	133	THR	N-CA-C	5.16	116.30	107.28
1	F	297	ALA	CB-CA-C	-5.13	109.68	117.07
1	F	253	ILE	N-CA-C	-5.13	100.17	107.77
1	A	173	SER	N-CA-C	-5.11	100.69	108.76
1	A	261	THR	N-CA-C	5.10	116.87	110.24
1	F	136	THR	N-CA-C	-5.06	107.10	113.28
1	G	290	GLN	N-CA-C	5.04	117.11	108.90
1	B	262	SER	N-CA-C	5.03	116.46	111.07
1	D	284	ASN	N-CA-C	5.03	118.25	112.57
1	D	209	SER	N-CA-C	5.02	117.58	110.50

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	1496	0	1453	159	1
1	B	1496	0	1453	159	0
1	D	1496	0	1453	166	0
1	E	1496	0	1453	156	1
1	F	1496	0	1453	137	0
1	G	1496	0	1453	164	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	8976	0	8718	876	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 50.

All (876) 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:E:217:ASN:ND2	1:E:224:SER:H	1.57	1.00
1:D:261:THR:HG22	1:D:301:PRO:HG3	1.45	0.98
1:A:143:ASN:HD21	1:A:157:LEU:H	1.09	0.96
1:B:268:PHE:HE1	1:B:300:SER:HA	1.27	0.96
1:F:242:PHE:CE1	1:F:277:ASN:ND2	2.38	0.92
1:F:225:GLU:HG3	1:F:226:THR:H	1.33	0.91
1:D:196:LEU:HA	1:D:209:SER:HB3	1.50	0.90
1:A:164:ASN:ND2	1:B:164:ASN:HD21	1.69	0.89
1:D:245:THR:HG22	1:D:246:THR:H	1.35	0.88
1:G:219:SER:HB2	1:G:224:SER:HB2	1.52	0.88
1:G:245:THR:HG22	1:G:246:THR:H	1.39	0.87
1:G:140:PRO:O	1:G:156:LYS:HE2	1.75	0.87
1:A:174:LEU:HD23	1:A:175:VAL:N	1.90	0.87
1:D:174:LEU:HD23	1:D:175:VAL:H	1.39	0.87
1:G:172:VAL:HG21	1:G:291:PHE:CE2	2.10	0.87
1:D:217:ASN:H	1:D:217:ASN:HD22	1.23	0.86
1:E:164:ASN:ND2	1:F:164:ASN:HD21	1.74	0.86
1:G:197:TYR:H	1:G:209:SER:HB3	1.41	0.85
1:D:217:ASN:ND2	1:D:224:SER:HB3	1.90	0.85
1:E:163:LYS:HE2	1:E:319:THR:HG21	1.59	0.85
1:E:217:ASN:ND2	1:E:217:ASN:H	1.72	0.85
1:D:270:LEU:HD23	1:D:270:LEU:H	1.39	0.85
1:D:174:LEU:HD23	1:D:175:VAL:N	1.91	0.85
1:A:164:ASN:HD21	1:D:164:ASN:ND2	1.75	0.84
1:B:225:GLU:HG3	1:B:226:THR:H	1.41	0.84
1:D:140:PRO:O	1:D:156:LYS:HE2	1.77	0.84
1:A:217:ASN:ND2	1:A:224:SER:H	1.77	0.83
1:D:282:SER:HB3	1:D:284:ASN:ND2	1.94	0.83
1:G:166:GLY:O	1:G:319:THR:HG22	1.77	0.83
1:E:219:SER:HB2	1:E:224:SER:HB2	1.61	0.83
1:E:193:GLN:HB3	1:E:292:GLU:HG3	1.59	0.82
1:B:142:PRO:HB3	1:B:177:VAL:HG21	1.60	0.82
1:G:217:ASN:H	1:G:217:ASN:HD22	1.24	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:217:ASN:HD21	1:F:224:SER:HB3	1.45	0.82
1:B:245:THR:HG22	1:B:246:THR:H	1.42	0.82
1:G:163:LYS:HE2	1:G:232:ALA:O	1.81	0.81
1:B:306:ALA:C	1:B:308:LEU:H	1.87	0.81
1:E:220:SER:HB3	1:E:222:ALA:H	1.46	0.81
1:F:268:PHE:HE1	1:F:300:SER:HA	1.44	0.81
1:G:225:GLU:HG3	1:G:226:THR:H	1.46	0.81
1:F:242:PHE:HE1	1:F:277:ASN:ND2	1.80	0.80
1:A:249:SER:CB	1:D:260:MET:HE1	2.11	0.80
1:F:245:THR:HG22	1:F:246:THR:H	1.45	0.80
1:A:216:LYS:HB2	1:A:223:THR:HB	1.64	0.80
1:E:245:THR:HG22	1:E:246:THR:H	1.45	0.80
1:D:225:GLU:HG3	1:D:226:THR:H	1.46	0.80
1:A:219:SER:HB2	1:A:224:SER:HB2	1.64	0.79
1:E:143:ASN:C	1:E:143:ASN:HD22	1.90	0.79
1:A:220:SER:HB3	1:A:222:ALA:H	1.48	0.79
1:G:261:THR:HG22	1:G:301:PRO:HG3	1.63	0.79
1:B:193:GLN:HB3	1:B:292:GLU:HG3	1.62	0.79
1:D:261:THR:HG22	1:D:301:PRO:CG	2.11	0.79
1:F:219:SER:HB2	1:F:224:SER:HB2	1.65	0.79
1:E:216:LYS:HB2	1:E:223:THR:HB	1.65	0.78
1:E:164:ASN:HD22	1:F:164:ASN:HD21	1.26	0.78
1:G:146:ILE:HD12	1:G:181:VAL:HG11	1.65	0.78
1:B:272:ILE:HA	1:B:292:GLU:O	1.84	0.78
1:E:143:ASN:HD21	1:E:157:LEU:H	1.30	0.78
1:A:219:SER:CB	1:A:224:SER:HB2	2.13	0.78
1:G:284:ASN:HD22	1:G:284:ASN:H	1.32	0.78
1:D:143:ASN:HA	1:D:153:ASN:HD22	1.49	0.77
1:A:193:GLN:HB3	1:A:292:GLU:HG3	1.66	0.77
1:G:172:VAL:HG23	1:G:315:PHE:HE2	1.48	0.77
1:E:217:ASN:HD21	1:E:224:SER:H	1.31	0.77
1:G:264:ASP:OD2	1:G:266:SER:HB2	1.84	0.77
1:F:217:ASN:C	1:F:217:ASN:HD22	1.91	0.77
1:G:131:ILE:C	1:G:131:ILE:HD12	2.09	0.77
1:G:268:PHE:HE1	1:G:300:SER:HA	1.49	0.77
1:A:164:ASN:HD21	1:D:164:ASN:HD21	1.32	0.77
1:B:268:PHE:CE1	1:B:300:SER:HA	2.18	0.77
1:D:298:SER:O	1:D:299:GLU:HB2	1.85	0.76
1:E:144:CYS:O	1:E:153:ASN:HA	1.85	0.76
1:D:219:SER:HB2	1:D:224:SER:HB2	1.67	0.76
1:D:277:ASN:HA	1:D:288:ALA:HB3	1.67	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:219:SER:CB	1:E:224:SER:HB2	2.16	0.76
1:E:298:SER:O	1:E:299:GLU:HB2	1.86	0.76
1:D:306:ALA:C	1:D:308:LEU:H	1.91	0.75
1:E:300:SER:OG	1:E:301:PRO:HD2	1.86	0.75
1:A:250:GLU:HB2	1:D:309:THR:HG21	1.68	0.75
1:E:259:TYR:HE2	1:E:301:PRO:HG2	1.50	0.75
1:G:298:SER:O	1:G:299:GLU:HB2	1.87	0.74
1:F:306:ALA:C	1:F:308:LEU:H	1.95	0.74
1:G:196:LEU:HA	1:G:209:SER:HB2	1.70	0.74
1:A:315:PHE:HA	1:D:314:PHE:CE1	2.23	0.74
1:B:247:ARG:O	1:B:250:GLU:HG2	1.88	0.74
1:D:166:GLY:O	1:D:319:THR:HG22	1.87	0.74
1:F:309:THR:HG21	1:G:250:GLU:HB2	1.67	0.74
1:A:205:LEU:HD21	1:A:284:ASN:O	1.87	0.74
1:B:264:ASP:OD2	1:B:266:SER:HB2	1.88	0.73
1:E:275:MET:O	1:E:276:LEU:HB2	1.87	0.73
1:E:315:PHE:HA	1:G:314:PHE:CE1	2.23	0.73
1:F:204:LEU:HD22	1:F:211:LEU:HD23	1.69	0.73
1:A:218:LYS:NZ	1:B:166:GLY:HA2	2.02	0.73
1:D:217:ASN:HD21	1:D:224:SER:HB3	1.51	0.73
1:F:264:ASP:OD2	1:F:266:SER:HB2	1.88	0.73
1:B:199:ASP:HB3	1:B:205:LEU:HD11	1.71	0.73
1:D:217:ASN:HD22	1:D:217:ASN:N	1.85	0.73
1:G:197:TYR:N	1:G:209:SER:HB3	2.04	0.73
1:E:227:VAL:C	1:E:229:SER:H	1.97	0.73
1:G:217:ASN:H	1:G:217:ASN:ND2	1.82	0.73
1:G:228:ALA:O	1:G:229:SER:HB3	1.88	0.73
1:B:251:ASN:C	1:B:251:ASN:HD22	1.96	0.72
1:D:245:THR:HG22	1:D:246:THR:N	2.04	0.72
1:F:308:LEU:HD23	1:F:308:LEU:O	1.89	0.72
1:E:164:ASN:HD21	1:G:164:ASN:HD21	1.36	0.72
1:F:244:THR:HG22	1:F:245:THR:O	1.89	0.72
1:A:258:TYR:C	1:A:270:LEU:HD21	2.15	0.72
1:B:143:ASN:HD21	1:B:157:LEU:H	1.33	0.72
1:F:193:GLN:HB3	1:F:292:GLU:HG3	1.72	0.72
1:G:144:CYS:SG	1:G:157:LEU:HB2	2.29	0.72
1:F:275:MET:O	1:F:275:MET:HE2	1.90	0.71
1:A:217:ASN:HD21	1:A:224:SER:H	1.38	0.71
1:D:217:ASN:H	1:D:217:ASN:ND2	1.87	0.71
1:D:304:ASN:C	1:D:306:ALA:H	1.95	0.71
1:B:182:ASN:ND2	1:B:307:THR:HA	2.05	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:304:ASN:C	1:D:306:ALA:N	2.46	0.71
1:A:298:SER:O	1:A:299:GLU:HB2	1.89	0.71
1:G:284:ASN:H	1:G:284:ASN:ND2	1.86	0.71
1:D:216:LYS:HB2	1:D:223:THR:HA	1.73	0.70
1:G:304:ASN:C	1:G:306:ALA:N	2.48	0.70
1:B:244:THR:HG22	1:B:245:THR:O	1.91	0.70
1:A:217:ASN:ND2	1:A:217:ASN:H	1.88	0.70
1:E:228:ALA:O	1:E:229:SER:HB3	1.92	0.70
1:A:300:SER:OG	1:A:301:PRO:HD2	1.90	0.70
1:D:159:LEU:HG	1:D:161:LEU:HD21	1.73	0.70
1:A:143:ASN:HD21	1:A:157:LEU:N	1.87	0.70
1:A:239:ALA:HB2	1:D:138:ILE:HG23	1.73	0.70
1:E:164:ASN:HD21	1:G:164:ASN:ND2	1.90	0.70
1:G:304:ASN:C	1:G:306:ALA:H	1.98	0.69
1:A:225:GLU:HG3	1:A:226:THR:H	1.57	0.69
1:G:216:LYS:HB2	1:G:223:THR:HA	1.73	0.69
1:A:164:ASN:ND2	1:D:164:ASN:HD21	1.90	0.69
1:E:172:VAL:HG22	1:E:315:PHE:HE2	1.58	0.69
1:G:261:THR:HG22	1:G:301:PRO:CG	2.22	0.69
1:G:219:SER:CB	1:G:224:SER:HB2	2.23	0.69
1:B:298:SER:O	1:B:299:GLU:HB2	1.93	0.69
1:E:183:GLN:HG2	1:E:186:THR:HG21	1.75	0.68
1:D:308:LEU:HD23	1:D:308:LEU:O	1.94	0.68
1:E:260:MET:HE1	1:F:249:SER:HB3	1.74	0.68
1:G:217:ASN:HD22	1:G:217:ASN:N	1.86	0.68
1:G:245:THR:HG22	1:G:246:THR:N	2.09	0.68
1:G:270:LEU:N	1:G:270:LEU:HD23	2.09	0.68
1:D:156:LYS:NZ	1:D:156:LYS:HB3	2.08	0.68
1:A:143:ASN:ND2	1:A:157:LEU:H	1.88	0.68
1:A:143:ASN:O	1:A:211:LEU:HD12	1.94	0.67
1:D:131:ILE:HD12	1:D:131:ILE:C	2.19	0.67
1:D:146:ILE:HD12	1:D:181:VAL:HG11	1.75	0.67
1:G:284:ASN:HD22	1:G:284:ASN:N	1.91	0.67
1:F:199:ASP:OD1	1:F:203:ASN:HB2	1.94	0.67
1:D:306:ALA:O	1:D:308:LEU:N	2.27	0.67
1:E:225:GLU:HG3	1:E:226:THR:H	1.58	0.67
1:B:251:ASN:O	1:B:251:ASN:ND2	2.27	0.67
1:E:172:VAL:HG11	1:E:291:PHE:CD2	2.30	0.67
1:B:304:ASN:C	1:B:306:ALA:N	2.49	0.67
1:A:183:GLN:HG2	1:A:186:THR:HG21	1.77	0.67
1:A:216:LYS:CB	1:A:223:THR:HB	2.24	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:143:ASN:HD21	1:F:157:LEU:H	1.41	0.67
1:E:250:GLU:HB2	1:G:309:THR:HG21	1.76	0.67
1:D:239:ALA:O	1:D:241:PRO:HD3	1.95	0.66
1:D:144:CYS:SG	1:D:157:LEU:HB2	2.35	0.66
1:B:306:ALA:C	1:B:308:LEU:N	2.50	0.66
1:D:261:THR:OG1	1:D:266:SER:N	2.28	0.66
1:D:161:LEU:HD22	1:D:161:LEU:N	2.11	0.66
1:A:144:CYS:O	1:A:153:ASN:HA	1.96	0.66
1:A:284:ASN:ND2	1:A:284:ASN:H	1.93	0.66
1:B:265:ARG:HE	1:B:265:ARG:HA	1.60	0.66
1:A:132:ASN:O	1:A:162:VAL:HA	1.96	0.65
1:A:195:ARG:HG2	1:A:290:GLN:HG3	1.77	0.65
1:E:253:ILE:HD13	1:G:314:PHE:HZ	1.61	0.65
1:G:277:ASN:HA	1:G:288:ALA:HB3	1.79	0.65
1:G:306:ALA:C	1:G:308:LEU:H	2.05	0.65
1:F:245:THR:HG22	1:F:246:THR:N	2.11	0.65
1:G:147:VAL:HB	1:G:150:THR:HG23	1.78	0.65
1:A:184:MET:HE2	1:A:190:ALA:HB1	1.79	0.65
1:E:287:TYR:HE2	1:E:320:GLU:OE1	1.78	0.65
1:E:131:ILE:C	1:E:131:ILE:HD12	2.22	0.65
1:E:136:THR:HG23	1:E:159:LEU:HB3	1.79	0.64
1:B:131:ILE:C	1:B:131:ILE:HD12	2.22	0.64
1:F:147:VAL:HG12	1:F:149:ASN:H	1.61	0.64
1:G:275:MET:SD	1:G:278:SER:HA	2.37	0.64
1:B:305:ILE:O	1:B:307:THR:N	2.30	0.64
1:G:239:ALA:O	1:G:241:PRO:HD3	1.98	0.64
1:A:249:SER:HB2	1:D:260:MET:HE1	1.79	0.64
1:E:217:ASN:H	1:E:217:ASN:HD22	1.46	0.64
1:G:270:LEU:HD23	1:G:270:LEU:H	1.63	0.64
1:E:132:ASN:O	1:E:162:VAL:HA	1.97	0.64
1:B:172:VAL:HG23	1:B:313:PHE:O	1.98	0.64
1:D:174:LEU:CD2	1:D:175:VAL:N	2.61	0.64
1:D:163:LYS:HE2	1:D:232:ALA:O	1.97	0.63
1:D:270:LEU:HD23	1:D:270:LEU:N	2.11	0.63
1:E:157:LEU:HD23	1:E:211:LEU:HD22	1.80	0.63
1:D:306:ALA:C	1:D:308:LEU:N	2.53	0.63
1:A:303:SER:O	1:A:304:ASN:HB2	1.98	0.63
1:F:150:THR:HB	1:F:152:THR:O	1.98	0.63
1:G:274:ILE:HA	1:G:290:GLN:O	1.97	0.63
1:E:164:ASN:ND2	1:G:164:ASN:HD21	1.94	0.63
1:F:228:ALA:O	1:F:229:SER:HB3	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:SER:C	1:A:251:ASN:H	2.07	0.62
1:B:172:VAL:HG22	1:B:315:PHE:HE2	1.63	0.62
1:E:172:VAL:HG22	1:E:315:PHE:CE2	2.34	0.62
1:A:304:ASN:O	1:A:306:ALA:N	2.31	0.62
1:A:131:ILE:HD12	1:A:131:ILE:C	2.24	0.62
1:B:138:ILE:HG12	1:D:318:ILE:HG21	1.80	0.62
1:B:245:THR:HG22	1:B:246:THR:N	2.14	0.62
1:F:131:ILE:HD12	1:F:131:ILE:C	2.24	0.62
1:D:205:LEU:HD23	1:D:208:GLU:OE1	1.99	0.62
1:F:225:GLU:CG	1:F:226:THR:H	2.06	0.62
1:A:172:VAL:HG22	1:A:315:PHE:HE2	1.64	0.62
1:G:303:SER:O	1:G:304:ASN:HB2	1.98	0.62
1:D:239:ALA:C	1:D:241:PRO:HD3	2.24	0.62
1:A:166:GLY:O	1:A:319:THR:HG22	2.00	0.62
1:B:207:GLU:OE2	1:B:207:GLU:HA	1.99	0.62
1:E:217:ASN:HD22	1:E:217:ASN:N	1.97	0.62
1:A:305:ILE:O	1:A:307:THR:N	2.33	0.61
1:E:249:SER:C	1:E:251:ASN:H	2.08	0.61
1:D:154:ASP:HB2	1:D:178:SER:HB3	1.81	0.61
1:D:282:SER:HB3	1:D:284:ASN:HD21	1.65	0.61
1:E:140:PRO:O	1:E:156:LYS:NZ	2.32	0.61
1:D:253:ILE:HG21	1:D:274:ILE:HD12	1.82	0.61
1:E:184:MET:C	1:E:186:THR:H	2.07	0.61
1:F:172:VAL:HG23	1:F:313:PHE:HB3	1.83	0.61
1:E:170:GLY:HA3	1:E:315:PHE:CZ	2.36	0.61
1:E:229:SER:OG	1:E:231:LYS:HG3	2.01	0.61
1:F:172:VAL:HG11	1:F:291:PHE:CE2	2.36	0.61
1:B:147:VAL:C	1:B:149:ASN:H	2.08	0.61
1:A:304:ASN:C	1:A:306:ALA:N	2.58	0.61
1:B:225:GLU:CG	1:B:226:THR:H	2.11	0.61
1:B:282:SER:HB3	1:B:284:ASN:ND2	2.15	0.60
1:B:300:SER:O	1:B:302:GLU:N	2.34	0.60
1:E:205:LEU:HD21	1:E:284:ASN:O	2.01	0.60
1:G:163:LYS:NZ	1:G:319:THR:HG21	2.15	0.60
1:A:167:LEU:HG	1:D:160:VAL:HG11	1.83	0.60
1:G:282:SER:HB3	1:G:284:ASN:ND2	2.16	0.60
1:A:228:ALA:O	1:A:229:SER:HB3	2.01	0.60
1:D:195:ARG:HG3	1:D:195:ARG:HH11	1.65	0.60
1:F:278:SER:OG	1:F:279:ARG:N	2.31	0.60
1:D:193:GLN:O	1:D:194:LEU:HD23	2.01	0.60
1:A:172:VAL:HG22	1:A:315:PHE:CE2	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:GLY:C	1:B:319:THR:HG22	2.27	0.60
1:E:304:ASN:C	1:E:306:ALA:N	2.57	0.60
1:G:282:SER:HB2	1:G:285:VAL:HG23	1.84	0.60
1:B:143:ASN:HD21	1:B:157:LEU:N	2.00	0.59
1:D:274:ILE:HA	1:D:290:GLN:O	2.02	0.59
1:G:131:ILE:HD12	1:G:131:ILE:O	2.01	0.59
1:G:195:ARG:HG2	1:G:290:GLN:NE2	2.17	0.59
1:G:196:LEU:HA	1:G:209:SER:CB	2.33	0.59
1:G:252:TYR:CE1	1:G:275:MET:HE3	2.37	0.59
1:A:164:ASN:HD22	1:B:164:ASN:HD21	1.49	0.59
1:E:302:GLU:O	1:E:302:GLU:HG3	2.01	0.59
1:B:141:PRO:O	1:B:143:ASN:N	2.36	0.59
1:B:227:VAL:O	1:F:227:VAL:HB	2.01	0.59
1:D:268:PHE:HE1	1:D:301:PRO:HD3	1.66	0.59
1:E:218:LYS:NZ	1:F:166:GLY:HA2	2.16	0.59
1:G:163:LYS:HE3	1:G:232:ALA:HB1	1.84	0.59
1:B:157:LEU:HD13	1:B:194:LEU:HD12	1.83	0.59
1:D:284:ASN:ND2	1:D:284:ASN:H	2.01	0.59
1:G:249:SER:C	1:G:251:ASN:H	2.10	0.59
1:A:164:ASN:HD21	1:B:164:ASN:HD21	1.49	0.59
1:A:251:ASN:H	1:A:251:ASN:HD22	1.50	0.59
1:D:159:LEU:HG	1:D:161:LEU:CD2	2.32	0.59
1:B:184:MET:HE2	1:B:190:ALA:HB1	1.85	0.59
1:D:228:ALA:O	1:D:229:SER:HB3	2.01	0.58
1:E:192:ILE:HG12	1:E:293:TRP:HB2	1.85	0.58
1:D:282:SER:HB2	1:D:285:VAL:HG23	1.84	0.58
1:G:308:LEU:C	1:G:308:LEU:HD23	2.28	0.58
1:A:218:LYS:HZ2	1:B:166:GLY:HA2	1.66	0.58
1:G:304:ASN:O	1:G:306:ALA:N	2.36	0.58
1:F:300:SER:O	1:F:302:GLU:N	2.36	0.58
1:B:302:GLU:OE1	1:B:302:GLU:C	2.47	0.58
1:G:253:ILE:O	1:G:273:SER:HA	2.03	0.58
1:F:270:LEU:HD12	1:F:293:TRP:HD1	1.67	0.58
1:A:259:TYR:HE1	1:A:308:LEU:HA	1.68	0.58
1:B:166:GLY:O	1:B:319:THR:HG22	2.04	0.58
1:B:182:ASN:HD22	1:B:307:THR:HA	1.68	0.58
1:F:275:MET:O	1:F:276:LEU:O	2.21	0.58
1:B:227:VAL:HB	1:F:227:VAL:O	2.03	0.58
1:D:193:GLN:HB3	1:D:292:GLU:HG3	1.85	0.58
1:G:239:ALA:C	1:G:241:PRO:HD3	2.28	0.58
1:G:251:ASN:H	1:G:251:ASN:HD22	1.51	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:251:ASN:C	1:B:251:ASN:ND2	2.61	0.58
1:E:272:ILE:HD13	1:E:293:TRP:NE1	2.18	0.58
1:F:147:VAL:C	1:F:149:ASN:H	2.12	0.58
1:E:147:VAL:C	1:E:149:ASN:H	2.12	0.57
1:E:242:PHE:HD2	1:E:279:ARG:HA	1.69	0.57
1:A:174:LEU:HD13	1:A:192:ILE:HD12	1.86	0.57
1:D:264:ASP:OD2	1:D:266:SER:HB2	2.04	0.57
1:F:160:VAL:HG21	1:G:318:ILE:HD11	1.86	0.57
1:G:261:THR:OG1	1:G:266:SER:N	2.37	0.57
1:G:132:ASN:O	1:G:162:VAL:HG13	2.04	0.57
1:B:242:PHE:O	1:B:243:ASN:C	2.46	0.57
1:E:225:GLU:CG	1:E:226:THR:H	2.17	0.57
1:A:218:LYS:HZ3	1:B:166:GLY:HA2	1.69	0.57
1:E:249:SER:HB2	1:G:260:MET:HE1	1.87	0.57
1:G:306:ALA:C	1:G:308:LEU:N	2.60	0.57
1:A:155:GLY:HA2	1:A:177:VAL:HG23	1.86	0.57
1:D:170:GLY:HA3	1:D:315:PHE:CE2	2.40	0.57
1:F:251:ASN:C	1:F:251:ASN:HD22	2.13	0.57
1:G:136:THR:CG2	1:G:159:LEU:HB3	2.34	0.57
1:G:168:VAL:HG23	1:G:319:THR:HB	1.87	0.57
1:G:261:THR:HA	1:G:301:PRO:HG2	1.86	0.57
1:F:172:VAL:O	1:F:312:PRO:HA	2.05	0.57
1:A:192:ILE:HG12	1:A:293:TRP:HB2	1.87	0.57
1:A:275:MET:O	1:A:276:LEU:HB2	2.05	0.57
1:F:163:LYS:HE3	1:F:232:ALA:HB1	1.87	0.57
1:F:300:SER:OG	1:F:301:PRO:HD2	2.05	0.57
1:B:217:ASN:HB3	1:B:224:SER:HB3	1.87	0.57
1:E:320:GLU:O	1:E:321:ASP:C	2.46	0.57
1:F:260:MET:HE1	1:G:249:SER:HB2	1.86	0.57
1:F:157:LEU:HD23	1:F:211:LEU:HD13	1.87	0.57
1:G:143:ASN:HD21	1:G:157:LEU:H	1.52	0.56
1:G:156:LYS:NZ	1:G:156:LYS:HB3	2.19	0.56
1:D:253:ILE:O	1:D:273:SER:HA	2.05	0.56
1:E:216:LYS:CB	1:E:223:THR:HB	2.33	0.56
1:A:164:ASN:ND2	1:D:164:ASN:ND2	2.48	0.56
1:A:245:THR:C	1:A:247:ARG:H	2.14	0.56
1:A:249:SER:O	1:A:251:ASN:N	2.39	0.56
1:B:308:LEU:HD23	1:B:308:LEU:O	2.06	0.56
1:D:143:ASN:CA	1:D:153:ASN:HD22	2.17	0.56
1:F:268:PHE:CE1	1:F:300:SER:HA	2.34	0.56
1:F:300:SER:O	1:F:301:PRO:C	2.49	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:147:VAL:HG12	1:B:149:ASN:H	1.70	0.56
1:E:216:LYS:HB2	1:E:223:THR:CB	2.34	0.56
1:G:217:ASN:ND2	1:G:224:SER:HB3	2.20	0.56
1:D:251:ASN:HD22	1:D:251:ASN:H	1.54	0.56
1:F:228:ALA:O	1:F:229:SER:CB	2.52	0.56
1:F:302:GLU:C	1:F:302:GLU:OE1	2.48	0.56
1:E:168:VAL:O	1:E:316:SER:HA	2.05	0.56
1:F:217:ASN:ND2	1:F:224:SER:HB3	2.18	0.56
1:E:167:LEU:HD23	1:E:318:ILE:HD12	1.87	0.56
1:D:300:SER:OG	1:D:301:PRO:HD2	2.06	0.56
1:D:245:THR:CG2	1:D:246:THR:H	2.14	0.56
1:E:261:THR:HG22	1:E:301:PRO:HG3	1.87	0.56
1:G:150:THR:HG21	1:G:154:ASP:OD2	2.06	0.56
1:B:196:LEU:HB2	1:B:289:ILE:O	2.06	0.55
1:B:141:PRO:O	1:B:142:PRO:C	2.50	0.55
1:D:143:ASN:HA	1:D:153:ASN:ND2	2.19	0.55
1:F:261:THR:HG22	1:F:301:PRO:CG	2.36	0.55
1:A:239:ALA:HB2	1:D:138:ILE:CG2	2.36	0.55
1:B:228:ALA:O	1:B:229:SER:HB3	2.05	0.55
1:E:298:SER:O	1:E:299:GLU:CB	2.55	0.55
1:F:138:ILE:HG12	1:G:318:ILE:HG21	1.89	0.55
1:A:259:TYR:HE2	1:A:301:PRO:HG2	1.69	0.55
1:B:189:THR:HG22	1:B:190:ALA:N	2.21	0.55
1:D:253:ILE:HB	1:D:274:ILE:HB	1.89	0.55
1:E:239:ALA:HB2	1:G:138:ILE:HG23	1.88	0.55
1:E:304:ASN:O	1:E:305:ILE:C	2.50	0.55
1:E:149:ASN:O	1:E:150:THR:O	2.25	0.55
1:F:135:TRP:CE2	1:F:137:GLY:HA2	2.40	0.55
1:G:300:SER:OG	1:G:301:PRO:HD2	2.06	0.55
1:E:167:LEU:CD2	1:E:318:ILE:HD12	2.37	0.55
1:F:237:THR:HG22	1:F:277:ASN:ND2	2.21	0.55
1:G:275:MET:HE2	1:G:275:MET:O	2.06	0.55
1:A:227:VAL:C	1:A:229:SER:H	2.12	0.55
1:A:307:THR:O	1:A:308:LEU:HB3	2.05	0.55
1:B:284:ASN:ND2	1:B:284:ASN:H	2.04	0.55
1:D:150:THR:HG22	1:D:152:THR:O	2.06	0.55
1:E:251:ASN:H	1:E:251:ASN:HD22	1.52	0.55
1:A:280:MET:HA	1:A:285:VAL:HG11	1.89	0.55
1:B:275:MET:O	1:B:276:LEU:O	2.25	0.55
1:A:147:VAL:C	1:A:149:ASN:H	2.15	0.55
1:B:158:THR:HB	1:B:173:SER:OG	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:140:PRO:O	1:B:156:LYS:NZ	2.39	0.55
1:D:168:VAL:HG23	1:D:319:THR:HB	1.88	0.55
1:G:193:GLN:HB3	1:G:292:GLU:HG3	1.87	0.55
1:D:179:ASP:HA	1:D:182:ASN:HB2	1.89	0.54
1:F:242:PHE:O	1:F:243:ASN:C	2.49	0.54
1:F:298:SER:O	1:F:299:GLU:HB2	2.07	0.54
1:B:304:ASN:C	1:B:306:ALA:H	2.15	0.54
1:A:280:MET:HB3	1:A:285:VAL:HB	1.88	0.54
1:A:320:GLU:O	1:A:321:ASP:C	2.51	0.54
1:B:159:LEU:HD12	1:B:171:TYR:O	2.06	0.54
1:D:275:MET:SD	1:D:278:SER:HA	2.47	0.54
1:E:135:TRP:CH2	1:E:138:ILE:HG13	2.42	0.54
1:E:258:TYR:C	1:E:270:LEU:HD21	2.32	0.54
1:E:304:ASN:O	1:E:306:ALA:N	2.41	0.54
1:F:204:LEU:HG	1:F:204:LEU:O	2.06	0.54
1:F:302:GLU:OE1	1:F:303:SER:N	2.40	0.54
1:E:251:ASN:O	1:E:275:MET:O	2.26	0.54
1:G:244:THR:HG22	1:G:245:THR:O	2.07	0.54
1:G:254:HIS:ND1	1:G:273:SER:HB3	2.23	0.54
1:E:174:LEU:HD12	1:E:293:TRP:CZ3	2.42	0.54
1:F:251:ASN:HD21	1:F:275:MET:CE	2.20	0.54
1:B:207:GLU:OE2	1:F:283:SER:HB3	2.07	0.54
1:E:172:VAL:HG11	1:E:291:PHE:CE2	2.43	0.54
1:E:249:SER:C	1:E:251:ASN:N	2.65	0.54
1:A:249:SER:HB3	1:D:260:MET:HE1	1.89	0.54
1:B:154:ASP:HB2	1:B:178:SER:HB3	1.89	0.54
1:F:141:PRO:O	1:F:143:ASN:N	2.40	0.54
1:G:298:SER:O	1:G:299:GLU:CB	2.56	0.54
1:D:305:ILE:O	1:D:307:THR:N	2.37	0.53
1:B:135:TRP:CE2	1:B:137:GLY:HA2	2.44	0.53
1:B:309:THR:HG21	1:D:250:GLU:HB2	1.90	0.53
1:G:265:ARG:HE	1:G:265:ARG:HA	1.72	0.53
1:A:282:SER:HB2	1:A:285:VAL:HG23	1.91	0.53
1:F:184:MET:C	1:F:186:THR:H	2.14	0.53
1:B:251:ASN:O	1:B:275:MET:O	2.26	0.53
1:D:219:SER:CB	1:D:224:SER:HB2	2.34	0.53
1:G:282:SER:CB	1:G:284:ASN:ND2	2.72	0.53
1:A:304:ASN:O	1:A:305:ILE:C	2.52	0.53
1:B:303:SER:OG	1:B:304:ASN:N	2.40	0.53
1:F:275:MET:O	1:F:276:LEU:C	2.52	0.53
1:G:172:VAL:HG23	1:G:315:PHE:CE2	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:261:THR:HB	1:A:301:PRO:HG3	1.90	0.52
1:D:268:PHE:CE1	1:D:301:PRO:HD3	2.45	0.52
1:E:143:ASN:O	1:E:211:LEU:HD12	2.09	0.52
1:E:263:TYR:H	1:E:263:TYR:HD1	1.56	0.52
1:F:304:ASN:C	1:F:306:ALA:N	2.60	0.52
1:G:282:SER:HB3	1:G:284:ASN:HD21	1.72	0.52
1:A:131:ILE:C	1:A:133:THR:H	2.16	0.52
1:E:196:LEU:HA	1:E:209:SER:HB3	1.91	0.52
1:E:260:MET:HE1	1:F:249:SER:CB	2.38	0.52
1:B:172:VAL:O	1:B:312:PRO:HA	2.09	0.52
1:B:174:LEU:HD23	1:B:175:VAL:N	2.25	0.52
1:E:220:SER:HB3	1:E:222:ALA:N	2.21	0.52
1:A:260:MET:HE1	1:B:249:SER:HB3	1.90	0.52
1:B:253:ILE:HB	1:B:274:ILE:HB	1.91	0.52
1:F:160:VAL:HB	1:F:171:TYR:HB3	1.91	0.52
1:F:306:ALA:C	1:F:308:LEU:N	2.58	0.52
1:A:242:PHE:CE2	1:A:280:MET:HG2	2.45	0.52
1:F:141:PRO:O	1:F:142:PRO:C	2.50	0.52
1:G:156:LYS:HG3	1:G:177:VAL:HG21	1.91	0.52
1:G:242:PHE:HD2	1:G:279:ARG:HA	1.73	0.52
1:A:198:PHE:HA	1:A:203:ASN:O	2.09	0.52
1:B:260:MET:HE3	1:B:265:ARG:CZ	2.40	0.52
1:D:156:LYS:NZ	1:D:156:LYS:CB	2.73	0.52
1:D:244:THR:HG22	1:D:245:THR:O	2.10	0.52
1:B:227:VAL:C	1:B:229:SER:H	2.16	0.52
1:B:249:SER:O	1:B:251:ASN:N	2.42	0.52
1:D:225:GLU:CG	1:D:226:THR:H	2.19	0.52
1:B:314:PHE:CE1	1:D:315:PHE:HA	2.45	0.52
1:E:132:ASN:O	1:E:162:VAL:HG13	2.10	0.52
1:F:131:ILE:HD12	1:F:131:ILE:O	2.10	0.52
1:A:318:ILE:O	1:A:318:ILE:HG22	2.09	0.51
1:E:225:GLU:HG3	1:E:226:THR:N	2.25	0.51
1:F:204:LEU:CD2	1:F:211:LEU:HD23	2.38	0.51
1:F:258:TYR:N	1:F:258:TYR:CD1	2.78	0.51
1:F:309:THR:HG21	1:G:250:GLU:CB	2.38	0.51
1:B:163:LYS:CE	1:B:232:ALA:HB1	2.40	0.51
1:G:131:ILE:C	1:G:131:ILE:CD1	2.77	0.51
1:A:272:ILE:HG13	1:A:273:SER:N	2.24	0.51
1:B:150:THR:HB	1:B:152:THR:O	2.09	0.51
1:B:257:CYS:C	1:B:258:TYR:CD1	2.88	0.51
1:D:231:LYS:HA	1:D:234:MET:HG2	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:ASN:HD21	1:B:247:ARG:NH2	2.07	0.51
1:A:143:ASN:CA	1:A:153:ASN:HD22	2.23	0.51
1:A:308:LEU:HD23	1:A:308:LEU:C	2.36	0.51
1:B:272:ILE:HD13	1:B:313:PHE:CD2	2.45	0.51
1:G:242:PHE:O	1:G:243:ASN:C	2.54	0.51
1:G:251:ASN:H	1:G:251:ASN:ND2	2.09	0.51
1:B:160:VAL:HB	1:B:171:TYR:HB3	1.91	0.51
1:E:163:LYS:HD3	1:E:232:ALA:O	2.10	0.51
1:E:227:VAL:O	1:E:229:SER:N	2.42	0.51
1:F:314:PHE:CE1	1:G:315:PHE:HA	2.46	0.51
1:A:245:THR:O	1:A:246:THR:OG1	2.21	0.51
1:A:270:LEU:HD23	1:A:270:LEU:N	2.25	0.51
1:D:143:ASN:C	1:D:143:ASN:HD22	2.18	0.51
1:D:146:ILE:HG22	1:D:146:ILE:O	2.10	0.51
1:D:284:ASN:H	1:D:284:ASN:HD22	1.59	0.51
1:G:261:THR:O	1:G:264:ASP:O	2.29	0.51
1:A:191:ASN:HD21	1:A:193:GLN:HE21	1.59	0.51
1:D:303:SER:O	1:D:304:ASN:HB2	2.10	0.51
1:G:189:THR:OG1	1:G:296:ASN:HA	2.10	0.51
1:A:284:ASN:N	1:A:284:ASN:HD22	2.09	0.51
1:E:227:VAL:C	1:E:229:SER:N	2.61	0.51
1:F:182:ASN:ND2	1:F:307:THR:HA	2.26	0.51
1:F:304:ASN:O	1:F:306:ALA:N	2.44	0.51
1:B:260:MET:HA	1:B:266:SER:O	2.11	0.50
1:D:197:TYR:N	1:D:209:SER:OG	2.39	0.50
1:F:226:THR:HG22	1:F:228:ALA:HB2	1.93	0.50
1:D:131:ILE:C	1:D:131:ILE:CD1	2.85	0.50
1:G:217:ASN:ND2	1:G:217:ASN:N	2.47	0.50
1:G:245:THR:CG2	1:G:246:THR:H	2.19	0.50
1:B:219:SER:HB2	1:B:224:SER:HB2	1.93	0.50
1:B:226:THR:HG22	1:B:228:ALA:HB2	1.93	0.50
1:D:170:GLY:HA3	1:D:315:PHE:CZ	2.47	0.50
1:F:135:TRP:CH2	1:F:218:LYS:HD2	2.47	0.50
1:F:284:ASN:ND2	1:F:284:ASN:H	2.09	0.50
1:G:245:THR:C	1:G:247:ARG:N	2.69	0.50
1:B:193:GLN:NE2	1:B:292:GLU:OE2	2.45	0.50
1:A:259:TYR:CE1	1:A:308:LEU:HA	2.46	0.50
1:B:244:THR:HG22	1:B:245:THR:N	2.27	0.50
1:G:143:ASN:C	1:G:143:ASN:HD22	2.19	0.50
1:A:154:ASP:C	1:A:177:VAL:HB	2.36	0.50
1:A:304:ASN:C	1:A:306:ALA:H	2.19	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:199:ASP:HB3	1:D:205:LEU:HD11	1.93	0.50
1:E:143:ASN:ND2	1:E:157:LEU:H	2.05	0.50
1:G:161:LEU:HD12	1:G:168:VAL:HG11	1.93	0.50
1:G:249:SER:O	1:G:251:ASN:N	2.45	0.50
1:A:143:ASN:C	1:A:143:ASN:HD22	2.20	0.50
1:A:287:TYR:HE2	1:A:320:GLU:OE1	1.95	0.50
1:B:277:ASN:O	1:B:278:SER:O	2.29	0.50
1:E:183:GLN:HG2	1:E:186:THR:CG2	2.41	0.50
1:A:170:GLY:HA3	1:A:315:PHE:CZ	2.46	0.49
1:A:171:TYR:CE2	1:B:318:ILE:HD11	2.47	0.49
1:D:257:CYS:O	1:D:270:LEU:HD23	2.12	0.49
1:E:252:TYR:CE2	1:E:275:MET:HG2	2.47	0.49
1:E:261:THR:HG22	1:E:301:PRO:CG	2.40	0.49
1:A:302:GLU:HG3	1:A:302:GLU:O	2.12	0.49
1:B:132:ASN:N	1:B:132:ASN:HD22	2.10	0.49
1:B:168:VAL:HG23	1:B:319:THR:HB	1.95	0.49
1:D:147:VAL:HB	1:D:150:THR:OG1	2.11	0.49
1:D:275:MET:O	1:D:276:LEU:HB2	2.11	0.49
1:D:307:THR:HG23	1:D:307:THR:O	2.11	0.49
1:F:225:GLU:HG3	1:F:226:THR:N	2.15	0.49
1:B:261:THR:HG22	1:B:301:PRO:HG3	1.94	0.49
1:D:132:ASN:O	1:D:162:VAL:HG13	2.12	0.49
1:E:167:LEU:HG	1:G:160:VAL:HG11	1.95	0.49
1:G:226:THR:HG22	1:G:228:ALA:HB2	1.93	0.49
1:A:141:PRO:O	1:A:142:PRO:C	2.55	0.49
1:A:198:PHE:CE2	1:A:204:LEU:HD13	2.47	0.49
1:A:227:VAL:C	1:A:229:SER:N	2.71	0.49
1:B:131:ILE:HD12	1:B:131:ILE:O	2.12	0.49
1:B:304:ASN:O	1:B:305:ILE:C	2.56	0.49
1:D:220:SER:HB3	1:D:222:ALA:H	1.77	0.49
1:D:263:TYR:CG	1:D:264:ASP:N	2.77	0.49
1:F:304:ASN:O	1:F:305:ILE:C	2.55	0.49
1:D:156:LYS:HB3	1:D:156:LYS:HZ3	1.74	0.49
1:D:305:ILE:O	1:D:305:ILE:HG22	2.12	0.49
1:G:270:LEU:N	1:G:270:LEU:CD2	2.75	0.49
1:B:144:CYS:SG	1:B:157:LEU:HB2	2.53	0.49
1:B:157:LEU:HD13	1:B:194:LEU:CD1	2.43	0.49
1:D:270:LEU:N	1:D:270:LEU:CD2	2.76	0.49
1:A:284:ASN:H	1:A:284:ASN:HD22	1.59	0.49
1:B:225:GLU:O	1:B:226:THR:HB	2.12	0.49
1:G:306:ALA:O	1:G:308:LEU:N	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:240:TYR:OH	1:D:318:ILE:N	2.42	0.49
1:A:225:GLU:CG	1:A:226:THR:H	2.25	0.48
1:E:282:SER:HB2	1:E:285:VAL:HG23	1.94	0.48
1:A:136:THR:HG23	1:A:159:LEU:HB3	1.94	0.48
1:A:184:MET:C	1:A:186:THR:H	2.21	0.48
1:D:149:ASN:O	1:D:150:THR:O	2.30	0.48
1:E:164:ASN:ND2	1:F:164:ASN:ND2	2.55	0.48
1:F:147:VAL:HG12	1:F:149:ASN:N	2.27	0.48
1:F:220:SER:HB3	1:F:222:ALA:H	1.78	0.48
1:A:216:LYS:HB2	1:A:223:THR:CB	2.41	0.48
1:E:141:PRO:O	1:E:142:PRO:C	2.57	0.48
1:G:141:PRO:O	1:G:142:PRO:C	2.55	0.48
1:G:302:GLU:OE1	1:G:303:SER:N	2.47	0.48
1:A:258:TYR:HA	1:A:270:LEU:CD2	2.43	0.48
1:D:176:GLY:H	1:D:308:LEU:HD23	1.79	0.48
1:E:149:ASN:O	1:E:150:THR:C	2.55	0.48
1:G:193:GLN:O	1:G:194:LEU:HD23	2.14	0.48
1:A:174:LEU:HD13	1:A:192:ILE:CD1	2.44	0.48
1:B:143:ASN:ND2	1:B:157:LEU:H	2.07	0.48
1:D:282:SER:CB	1:D:284:ASN:ND2	2.72	0.48
1:F:231:LYS:HE3	1:F:322:ASP:O	2.14	0.48
1:G:195:ARG:HG3	1:G:195:ARG:HH11	1.78	0.48
1:G:199:ASP:OD1	1:G:199:ASP:C	2.57	0.48
1:G:251:ASN:HD22	1:G:251:ASN:N	2.09	0.48
1:A:249:SER:C	1:A:251:ASN:N	2.70	0.48
1:E:252:TYR:CE1	1:E:275:MET:HE3	2.49	0.48
1:G:302:GLU:OE1	1:G:302:GLU:C	2.57	0.48
1:A:219:SER:O	1:A:220:SER:HB2	2.12	0.47
1:A:229:SER:OG	1:A:231:LYS:HG3	2.14	0.47
1:B:281:ILE:HG13	1:B:281:ILE:O	2.13	0.47
1:E:284:ASN:H	1:E:284:ASN:ND2	2.10	0.47
1:E:314:PHE:CE1	1:F:315:PHE:HA	2.48	0.47
1:B:181:VAL:C	1:B:183:GLN:H	2.22	0.47
1:D:144:CYS:O	1:D:153:ASN:HA	2.13	0.47
1:D:149:ASN:O	1:D:150:THR:C	2.57	0.47
1:D:277:ASN:O	1:D:278:SER:C	2.57	0.47
1:F:265:ARG:HE	1:F:265:ARG:HA	1.79	0.47
1:G:277:ASN:O	1:G:278:SER:C	2.57	0.47
1:B:162:VAL:HG11	1:D:164:ASN:O	2.13	0.47
1:B:163:LYS:HE3	1:B:232:ALA:HB1	1.95	0.47
1:E:219:SER:O	1:E:220:SER:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:307:THR:O	1:E:308:LEU:HB3	2.15	0.47
1:G:268:PHE:CE1	1:G:300:SER:HA	2.39	0.47
1:A:275:MET:HB2	1:A:290:GLN:HB3	1.96	0.47
1:B:303:SER:O	1:B:304:ASN:HB2	2.15	0.47
1:B:304:ASN:O	1:B:306:ALA:N	2.47	0.47
1:E:184:MET:C	1:E:186:THR:N	2.71	0.47
1:E:236:SER:HB2	1:E:318:ILE:HG22	1.96	0.47
1:A:174:LEU:HD23	1:A:175:VAL:H	1.76	0.47
1:D:163:LYS:HE3	1:D:232:ALA:HB1	1.96	0.47
1:G:156:LYS:HB3	1:G:156:LYS:HZ3	1.78	0.47
1:G:185:PHE:O	1:G:300:SER:OG	2.30	0.47
1:A:192:ILE:CG1	1:A:293:TRP:HB2	2.44	0.47
1:D:185:PHE:O	1:D:187:GLN:N	2.48	0.47
1:E:198:PHE:HB3	1:E:203:ASN:O	2.14	0.47
1:F:164:ASN:ND2	1:G:164:ASN:HD21	2.12	0.47
1:A:133:THR:O	1:A:217:ASN:HA	2.14	0.47
1:B:164:ASN:ND2	1:D:164:ASN:HD21	2.13	0.47
1:D:161:LEU:N	1:D:161:LEU:CD2	2.77	0.47
1:D:272:ILE:HD13	1:D:313:PHE:CD2	2.50	0.47
1:F:189:THR:HG22	1:F:190:ALA:N	2.29	0.47
1:F:274:ILE:HA	1:F:290:GLN:O	2.14	0.47
1:A:251:ASN:O	1:A:275:MET:O	2.32	0.47
1:F:245:THR:O	1:F:247:ARG:N	2.48	0.47
1:B:174:LEU:HD23	1:B:175:VAL:H	1.79	0.47
1:D:251:ASN:HD22	1:D:251:ASN:N	2.11	0.47
1:E:155:GLY:HA2	1:E:177:VAL:HG23	1.96	0.47
1:E:185:PHE:O	1:E:300:SER:HB3	2.15	0.47
1:E:220:SER:CB	1:E:222:ALA:H	2.23	0.47
1:B:282:SER:O	1:B:285:VAL:HG23	2.15	0.47
1:D:242:PHE:CG	1:D:280:MET:HE3	2.50	0.47
1:A:217:ASN:N	1:A:217:ASN:HD22	2.12	0.46
1:B:220:SER:HB3	1:B:222:ALA:H	1.80	0.46
1:B:261:THR:O	1:B:264:ASP:O	2.33	0.46
1:F:138:ILE:O	1:F:139:ASN:O	2.33	0.46
1:A:200:SER:O	1:A:231:LYS:HD3	2.16	0.46
1:A:263:TYR:CG	1:A:264:ASP:N	2.79	0.46
1:B:189:THR:CG2	1:B:190:ALA:N	2.78	0.46
1:D:150:THR:CG2	1:D:152:THR:O	2.63	0.46
1:F:144:CYS:HB3	1:F:145:GLN:H	1.50	0.46
1:G:170:GLY:HA3	1:G:315:PHE:CE2	2.50	0.46
1:A:159:LEU:HD23	1:A:215:LEU:HD21	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:135:TRP:CD1	1:D:135:TRP:C	2.93	0.46
1:D:258:TYR:HA	1:D:270:LEU:CD2	2.45	0.46
1:E:249:SER:CB	1:G:260:MET:HE1	2.45	0.46
1:F:135:TRP:CD1	1:F:135:TRP:C	2.92	0.46
1:E:131:ILE:C	1:E:131:ILE:CD1	2.89	0.46
1:F:219:SER:O	1:F:220:SER:HB2	2.16	0.46
1:G:258:TYR:HA	1:G:270:LEU:CD2	2.45	0.46
1:B:199:ASP:CB	1:B:205:LEU:HD11	2.44	0.46
1:D:198:PHE:C	1:D:205:LEU:HD13	2.41	0.46
1:E:195:ARG:HG2	1:E:290:GLN:HG3	1.98	0.46
1:F:258:TYR:C	1:F:270:LEU:HD21	2.40	0.46
1:B:184:MET:C	1:B:186:THR:H	2.21	0.46
1:E:166:GLY:O	1:E:319:THR:HG22	2.16	0.46
1:E:174:LEU:HD23	1:E:175:VAL:H	1.81	0.46
1:G:263:TYR:CG	1:G:264:ASP:N	2.84	0.46
1:A:252:TYR:CE2	1:A:275:MET:HG2	2.51	0.46
1:F:204:LEU:HD11	1:F:209:SER:OG	2.15	0.46
1:B:147:VAL:HG12	1:B:149:ASN:N	2.31	0.46
1:D:156:LYS:HG3	1:D:177:VAL:CG2	2.46	0.46
1:G:225:GLU:CG	1:G:226:THR:H	2.18	0.46
1:A:191:ASN:ND2	1:A:193:GLN:HE21	2.14	0.45
1:A:318:ILE:HD12	1:A:318:ILE:N	2.31	0.45
1:F:162:VAL:HG11	1:G:164:ASN:O	2.16	0.45
1:F:217:ASN:C	1:F:217:ASN:ND2	2.64	0.45
1:D:154:ASP:CB	1:D:178:SER:HB3	2.46	0.45
1:F:270:LEU:HD12	1:F:293:TRP:CD1	2.50	0.45
1:A:306:ALA:C	1:A:308:LEU:N	2.73	0.45
1:B:135:TRP:CZ3	1:B:138:ILE:HG13	2.52	0.45
1:B:261:THR:OG1	1:B:264:ASP:O	2.33	0.45
1:E:162:VAL:HG11	1:F:164:ASN:OD1	2.16	0.45
1:F:253:ILE:HB	1:F:274:ILE:HB	1.98	0.45
1:G:280:MET:HE2	1:G:285:VAL:O	2.16	0.45
1:A:172:VAL:O	1:A:312:PRO:HA	2.17	0.45
1:A:303:SER:O	1:A:304:ASN:CB	2.65	0.45
1:D:304:ASN:O	1:D:306:ALA:N	2.48	0.45
1:E:251:ASN:H	1:E:251:ASN:ND2	2.15	0.45
1:F:245:THR:CG2	1:F:246:THR:H	2.22	0.45
1:F:272:ILE:HA	1:F:292:GLU:O	2.17	0.45
1:A:143:ASN:C	1:A:153:ASN:HD22	2.24	0.45
1:A:315:PHE:HA	1:D:314:PHE:CD1	2.50	0.45
1:E:280:MET:HA	1:E:285:VAL:HG11	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:MET:C	1:A:282:SER:H	2.24	0.45
1:D:150:THR:HG22	1:D:152:THR:H	1.82	0.45
1:D:300:SER:OG	1:D:301:PRO:CD	2.65	0.45
1:E:170:GLY:HA3	1:E:315:PHE:CE1	2.52	0.45
1:E:218:LYS:HZ2	1:F:166:GLY:HA2	1.79	0.45
1:E:249:SER:HB3	1:E:252:TYR:CE1	2.52	0.45
1:F:143:ASN:C	1:F:143:ASN:HD22	2.25	0.45
1:B:260:MET:HE3	1:B:265:ARG:NH2	2.32	0.45
1:D:178:SER:O	1:D:181:VAL:N	2.50	0.45
1:D:307:THR:O	1:D:308:LEU:O	2.34	0.45
1:G:197:TYR:H	1:G:209:SER:CB	2.20	0.45
1:G:252:TYR:CZ	1:G:275:MET:HE3	2.51	0.45
1:G:289:ILE:HD11	1:G:317:TYR:OH	2.17	0.45
1:F:242:PHE:CD2	1:F:280:MET:HG2	2.51	0.45
1:G:131:ILE:O	1:G:131:ILE:CD1	2.65	0.45
1:A:174:LEU:CD1	1:A:192:ILE:HD12	2.47	0.45
1:A:251:ASN:HD22	1:A:251:ASN:N	2.10	0.45
1:E:146:ILE:HD11	1:E:174:LEU:HD21	1.98	0.45
1:E:196:LEU:HA	1:E:209:SER:CB	2.47	0.45
1:E:275:MET:HB2	1:E:290:GLN:HB3	1.98	0.45
1:A:140:PRO:O	1:A:156:LYS:HE2	2.16	0.45
1:A:143:ASN:HA	1:A:153:ASN:ND2	2.32	0.45
1:F:261:THR:OG1	1:F:264:ASP:O	2.33	0.45
1:F:305:ILE:O	1:F:307:THR:N	2.50	0.45
1:A:225:GLU:HG3	1:A:226:THR:N	2.30	0.44
1:G:249:SER:C	1:G:251:ASN:N	2.73	0.44
1:A:131:ILE:O	1:A:133:THR:N	2.51	0.44
1:B:199:ASP:OD1	1:B:203:ASN:HB2	2.17	0.44
1:B:284:ASN:ND2	1:B:284:ASN:N	2.65	0.44
1:D:196:LEU:HB2	1:D:289:ILE:O	2.18	0.44
1:E:225:GLU:CG	1:E:226:THR:N	2.80	0.44
1:F:304:ASN:C	1:F:306:ALA:H	2.25	0.44
1:G:144:CYS:HG	1:G:157:LEU:HB2	1.82	0.44
1:B:149:ASN:O	1:B:150:THR:C	2.60	0.44
1:B:225:GLU:HG3	1:B:226:THR:N	2.21	0.44
1:B:320:GLU:O	1:B:321:ASP:C	2.59	0.44
1:D:185:PHE:O	1:D:300:SER:OG	2.35	0.44
1:E:142:PRO:HA	1:E:155:GLY:O	2.17	0.44
1:E:218:LYS:HZ3	1:F:166:GLY:HA2	1.83	0.44
1:E:251:ASN:HD22	1:E:251:ASN:N	2.12	0.44
1:E:257:CYS:HG	1:E:293:TRP:CD1	2.35	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:244:THR:HG22	1:A:245:THR:N	2.33	0.44
1:B:164:ASN:ND2	1:D:164:ASN:OD1	2.50	0.44
1:B:247:ARG:O	1:B:248:ASP:C	2.61	0.44
1:D:275:MET:O	1:D:275:MET:HE2	2.18	0.44
1:E:135:TRP:CD1	1:E:135:TRP:C	2.95	0.44
1:E:135:TRP:HB2	1:E:160:VAL:HG22	1.99	0.44
1:B:164:ASN:HD21	1:D:164:ASN:HD21	1.65	0.44
1:E:245:THR:HG22	1:E:246:THR:N	2.22	0.44
1:A:146:ILE:HB	1:A:154:ASP:OD1	2.18	0.44
1:G:217:ASN:HD21	1:G:224:SER:HB3	1.82	0.44
1:E:257:CYS:SG	1:E:293:TRP:HD1	2.41	0.44
1:F:261:THR:O	1:F:264:ASP:O	2.36	0.44
1:G:149:ASN:O	1:G:150:THR:C	2.60	0.44
1:A:163:LYS:HE2	1:A:319:THR:HG21	1.99	0.44
1:B:140:PRO:HB2	1:B:141:PRO:HD2	2.00	0.44
1:B:188:LYS:HD3	1:B:188:LYS:N	2.33	0.44
1:D:156:LYS:HG3	1:D:177:VAL:HG21	1.99	0.44
1:E:280:MET:C	1:E:282:SER:H	2.25	0.44
1:B:147:VAL:C	1:B:149:ASN:N	2.75	0.43
1:E:143:ASN:C	1:E:143:ASN:ND2	2.64	0.43
1:E:198:PHE:HA	1:E:203:ASN:O	2.17	0.43
1:F:196:LEU:HD11	1:F:211:LEU:HB2	1.98	0.43
1:G:282:SER:CB	1:G:284:ASN:HD21	2.29	0.43
1:G:305:ILE:O	1:G:307:THR:N	2.47	0.43
1:E:184:MET:HE2	1:E:190:ALA:HB1	2.00	0.43
1:G:144:CYS:O	1:G:153:ASN:HA	2.18	0.43
1:G:161:LEU:N	1:G:161:LEU:HD22	2.33	0.43
1:A:131:ILE:C	1:A:131:ILE:CD1	2.91	0.43
1:E:135:TRP:CE2	1:E:137:GLY:HA2	2.53	0.43
1:E:185:PHE:CE2	1:E:308:LEU:HG	2.53	0.43
1:F:219:SER:CB	1:F:224:SER:HB2	2.41	0.43
1:E:155:GLY:CA	1:E:177:VAL:HG23	2.48	0.43
1:F:303:SER:O	1:F:304:ASN:HB2	2.18	0.43
1:G:305:ILE:O	1:G:305:ILE:HG22	2.19	0.43
1:A:143:ASN:HA	1:A:153:ASN:HD22	1.81	0.43
1:A:300:SER:OG	1:A:301:PRO:CD	2.62	0.43
1:D:260:MET:HA	1:D:266:SER:O	2.19	0.43
1:E:131:ILE:C	1:E:133:THR:H	2.26	0.43
1:F:144:CYS:SG	1:F:157:LEU:HB2	2.58	0.43
1:F:245:THR:CG2	1:F:246:THR:N	2.81	0.43
1:F:284:ASN:H	1:F:284:ASN:HD22	1.65	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:156:LYS:HG3	1:G:177:VAL:CG2	2.47	0.43
1:G:172:VAL:HG21	1:G:291:PHE:CZ	2.51	0.43
1:D:284:ASN:HD22	1:D:284:ASN:N	2.15	0.43
1:E:280:MET:HE2	1:E:285:VAL:O	2.19	0.43
1:F:199:ASP:OD1	1:F:203:ASN:N	2.49	0.43
1:F:296:ASN:OD1	1:F:296:ASN:O	2.36	0.43
1:G:247:ARG:O	1:G:248:ASP:C	2.60	0.43
1:A:284:ASN:ND2	1:A:284:ASN:N	2.56	0.43
1:D:264:ASP:OD2	1:D:266:SER:CB	2.66	0.43
1:F:149:ASN:O	1:F:150:THR:C	2.61	0.43
1:F:191:ASN:OD1	1:F:294:ASN:OD1	2.37	0.43
1:A:140:PRO:O	1:A:156:LYS:NZ	2.47	0.43
1:A:248:ASP:O	1:A:249:SER:HB2	2.19	0.43
1:B:314:PHE:N	1:B:314:PHE:CD2	2.81	0.43
1:A:225:GLU:O	1:A:226:THR:HB	2.19	0.43
1:B:219:SER:O	1:B:220:SER:HB2	2.19	0.43
1:B:258:TYR:CD1	1:B:258:TYR:N	2.86	0.43
1:B:283:SER:HB3	1:F:207:GLU:OE2	2.18	0.43
1:D:192:ILE:HG13	1:D:293:TRP:HB2	2.00	0.43
1:B:174:LEU:CD2	1:B:175:VAL:N	2.82	0.43
1:B:219:SER:CB	1:B:224:SER:HB2	2.48	0.43
1:B:249:SER:HB2	1:B:250:GLU:H	1.66	0.43
1:D:143:ASN:C	1:D:153:ASN:HD22	2.27	0.43
1:D:245:THR:O	1:D:247:ARG:N	2.52	0.43
1:A:131:ILE:C	1:A:133:THR:N	2.76	0.42
1:D:146:ILE:CD1	1:D:181:VAL:HG11	2.47	0.42
1:D:164:ASN:HD22	1:D:169:ASN:ND2	2.17	0.42
1:E:147:VAL:O	1:E:149:ASN:N	2.52	0.42
1:E:272:ILE:HG13	1:E:273:SER:N	2.33	0.42
1:F:143:ASN:H	1:F:143:ASN:ND2	2.17	0.42
1:F:207:GLU:HB3	1:F:208:GLU:OE1	2.20	0.42
1:G:242:PHE:CD1	1:G:280:MET:HE3	2.54	0.42
1:E:253:ILE:O	1:E:273:SER:HA	2.17	0.42
1:D:295:LEU:HD12	1:D:296:ASN:H	1.84	0.42
1:F:140:PRO:HB2	1:F:141:PRO:HD2	2.01	0.42
1:A:252:TYR:HB3	1:A:254:HIS:CE1	2.55	0.42
1:G:304:ASN:O	1:G:305:ILE:C	2.62	0.42
1:A:168:VAL:CG2	1:A:319:THR:HB	2.50	0.42
1:B:231:LYS:HB3	1:B:231:LYS:HE3	1.81	0.42
1:B:234:MET:O	1:B:320:GLU:N	2.46	0.42
1:B:256:ILE:HG23	1:B:270:LEU:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:306:ALA:O	1:B:308:LEU:N	2.52	0.42
1:F:263:TYR:CG	1:F:264:ASP:N	2.88	0.42
1:B:156:LYS:HG3	1:B:177:VAL:CG2	2.49	0.42
1:E:168:VAL:HG23	1:E:319:THR:HB	2.01	0.42
1:F:320:GLU:O	1:F:321:ASP:C	2.62	0.42
1:G:156:LYS:NZ	1:G:156:LYS:CB	2.82	0.42
1:G:220:SER:HB3	1:G:222:ALA:H	1.84	0.42
1:A:261:THR:HG22	1:A:301:PRO:CG	2.49	0.42
1:B:244:THR:CG2	1:B:245:THR:N	2.82	0.42
1:E:141:PRO:HB2	1:E:142:PRO:HD2	2.01	0.42
1:F:257:CYS:C	1:F:258:TYR:CD1	2.97	0.42
1:G:228:ALA:O	1:G:229:SER:CB	2.62	0.42
1:A:253:ILE:O	1:A:273:SER:HA	2.20	0.42
1:B:131:ILE:HD12	1:B:132:ASN:HD22	1.85	0.42
1:A:167:LEU:HG	1:D:160:VAL:CG1	2.49	0.42
1:B:245:THR:CG2	1:B:246:THR:H	2.23	0.42
1:B:284:ASN:H	1:B:284:ASN:HD22	1.68	0.42
1:F:132:ASN:N	1:F:132:ASN:HD22	2.17	0.42
1:A:258:TYR:HA	1:A:270:LEU:HD23	2.02	0.42
1:A:275:MET:O	1:A:276:LEU:CB	2.67	0.42
1:B:135:TRP:CH2	1:B:138:ILE:HG13	2.54	0.42
1:D:195:ARG:HG3	1:D:195:ARG:NH1	2.33	0.42
1:D:240:TYR:CD1	1:D:276:LEU:HD13	2.54	0.42
1:E:147:VAL:C	1:E:149:ASN:N	2.77	0.42
1:E:248:ASP:CG	1:G:265:ARG:HH12	2.28	0.42
1:G:303:SER:O	1:G:304:ASN:CB	2.66	0.42
1:A:205:LEU:HD12	1:A:205:LEU:N	2.35	0.41
1:B:132:ASN:N	1:B:132:ASN:ND2	2.68	0.41
1:B:217:ASN:O	1:B:223:THR:HA	2.20	0.41
1:B:302:GLU:O	1:B:302:GLU:HG3	2.20	0.41
1:D:176:GLY:H	1:D:308:LEU:CD2	2.33	0.41
1:D:303:SER:OG	1:D:304:ASN:N	2.53	0.41
1:F:251:ASN:O	1:F:251:ASN:ND2	2.53	0.41
1:F:140:PRO:O	1:F:156:LYS:HE2	2.20	0.41
1:F:221:THR:O	1:F:221:THR:HG22	2.20	0.41
1:G:170:GLY:HA3	1:G:315:PHE:CZ	2.56	0.41
1:G:174:LEU:HD23	1:G:175:VAL:H	1.84	0.41
1:B:282:SER:HB3	1:B:284:ASN:HD21	1.82	0.41
1:F:251:ASN:HD21	1:F:275:MET:HE3	1.84	0.41
1:A:213:ILE:HA	1:A:214:PRO:HD3	1.95	0.41
1:E:304:ASN:C	1:E:306:ALA:H	2.26	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:223:THR:OG1	1:B:224:SER:N	2.53	0.41
1:D:261:THR:O	1:D:264:ASP:O	2.39	0.41
1:E:138:ILE:O	1:E:139:ASN:C	2.64	0.41
1:F:277:ASN:O	1:F:278:SER:O	2.38	0.41
1:G:205:LEU:HD21	1:G:284:ASN:O	2.20	0.41
1:A:265:ARG:NH1	1:B:249:SER:OG	2.53	0.41
1:B:197:TYR:HB3	1:B:285:VAL:HG13	2.02	0.41
1:D:248:ASP:O	1:D:250:GLU:HG2	2.21	0.41
1:E:133:THR:HB	1:E:218:LYS:HB3	2.03	0.41
1:E:157:LEU:CD2	1:E:211:LEU:HD22	2.48	0.41
1:E:300:SER:O	1:E:301:PRO:C	2.63	0.41
1:G:163:LYS:HZ1	1:G:319:THR:HG21	1.84	0.41
1:G:207:GLU:OE2	1:G:207:GLU:HA	2.21	0.41
1:G:275:MET:O	1:G:276:LEU:HB2	2.20	0.41
1:A:168:VAL:HG23	1:A:319:THR:HB	2.03	0.41
1:A:231:LYS:HD2	1:A:287:TYR:OH	2.21	0.41
1:D:147:VAL:HG12	1:D:148:GLU:N	2.36	0.41
1:D:183:GLN:O	1:D:184:MET:C	2.63	0.41
1:D:228:ALA:O	1:D:229:SER:CB	2.68	0.41
1:E:172:VAL:CG1	1:E:291:PHE:CE2	3.04	0.41
1:E:305:ILE:O	1:E:306:ALA:C	2.64	0.41
1:G:171:TYR:C	1:G:171:TYR:CD2	2.98	0.41
1:A:155:GLY:CA	1:A:177:VAL:HG23	2.51	0.41
1:A:259:TYR:N	1:A:270:LEU:HD21	2.35	0.41
1:B:143:ASN:C	1:B:143:ASN:HD22	2.29	0.41
1:B:172:VAL:HG22	1:B:315:PHE:CE2	2.48	0.41
1:B:206:THR:HG22	1:B:212:LYS:HA	2.03	0.41
1:B:231:LYS:HE3	1:B:322:ASP:O	2.20	0.41
1:D:183:GLN:O	1:D:186:THR:HG22	2.21	0.41
1:D:231:LYS:HA	1:D:234:MET:CG	2.51	0.41
1:D:248:ASP:O	1:D:249:SER:C	2.64	0.41
1:E:219:SER:OG	1:E:224:SER:HB2	2.20	0.41
1:F:205:LEU:HA	1:F:205:LEU:HD12	1.74	0.41
1:G:174:LEU:CD2	1:G:175:VAL:N	2.84	0.41
1:G:236:SER:C	1:G:238:THR:H	2.28	0.41
1:G:245:THR:CG2	1:G:246:THR:N	2.80	0.41
1:G:245:THR:C	1:G:247:ARG:H	2.29	0.41
1:A:164:ASN:HD21	1:D:164:ASN:CG	2.29	0.41
1:A:198:PHE:HA	1:A:205:LEU:HD13	2.03	0.41
1:A:261:THR:CB	1:A:301:PRO:HG3	2.50	0.41
1:G:199:ASP:HB3	1:G:205:LEU:HD11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:216:LYS:HB3	1:G:223:THR:HB	2.03	0.41
1:G:300:SER:OG	1:G:301:PRO:CD	2.68	0.41
1:E:203:ASN:N	1:E:203:ASN:HD22	2.19	0.40
1:E:257:CYS:SG	1:E:293:TRP:CD1	3.14	0.40
1:F:242:PHE:HD2	1:F:279:ARG:HA	1.86	0.40
1:F:280:MET:HB3	1:F:285:VAL:HB	2.02	0.40
1:G:143:ASN:HD22	1:G:144:CYS:N	2.19	0.40
1:B:282:SER:HB2	1:B:285:VAL:HG23	2.03	0.40
1:A:147:VAL:C	1:A:149:ASN:N	2.80	0.40
1:D:243:ASN:OD1	1:D:243:ASN:O	2.38	0.40
1:E:257:CYS:C	1:E:258:TYR:CD1	3.00	0.40
1:F:168:VAL:HG23	1:F:319:THR:HB	2.03	0.40
1:G:147:VAL:C	1:G:149:ASN:H	2.30	0.40
1:A:174:LEU:HD23	1:A:174:LEU:C	2.43	0.40
1:B:252:TYR:CE2	1:B:275:MET:HG3	2.56	0.40
1:D:216:LYS:HD3	1:D:223:THR:HB	2.03	0.40
1:D:247:ARG:O	1:D:248:ASP:C	2.65	0.40
1:D:257:CYS:SG	1:D:270:LEU:HG	2.61	0.40
1:E:134:LEU:HA	1:E:134:LEU:HD12	1.87	0.40
1:E:217:ASN:HD21	1:E:224:SER:N	2.07	0.40
1:G:135:TRP:HB2	1:G:136:THR:H	1.67	0.40
1:B:191:ASN:OD1	1:B:294:ASN:OD1	2.40	0.40
1:B:235:PRO:HG2	1:B:277:ASN:OD1	2.21	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:186:THR:O	1:E:302:GLU:OE1[7_444]	2.08	0.12

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	190/220 (86%)	139 (73%)	29 (15%)	22 (12%)	0	0
1	B	190/220 (86%)	132 (70%)	31 (16%)	27 (14%)	0	0
1	D	190/220 (86%)	139 (73%)	30 (16%)	21 (11%)	0	0
1	E	190/220 (86%)	135 (71%)	32 (17%)	23 (12%)	0	0
1	F	190/220 (86%)	139 (73%)	28 (15%)	23 (12%)	0	0
1	G	190/220 (86%)	136 (72%)	30 (16%)	24 (13%)	0	0
All	All	1140/1320 (86%)	820 (72%)	180 (16%)	140 (12%)	0	0

All (140) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	220	SER
1	A	221	THR
1	A	248	ASP
1	A	250	GLU
1	A	304	ASN
1	A	306	ALA
1	B	145	GLN
1	B	220	SER
1	B	221	THR
1	B	223	THR
1	B	250	GLU
1	B	276	LEU
1	B	278	SER
1	B	301	PRO
1	B	304	ASN
1	B	306	ALA
1	D	150	THR
1	D	221	THR
1	D	243	ASN
1	D	262	SER
1	D	299	GLU
1	D	304	ASN
1	D	306	ALA
1	D	308	LEU
1	E	150	THR
1	E	220	SER
1	E	221	THR
1	E	248	ASP
1	E	250	GLU
1	E	299	GLU

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Mol	Chain	Res	Type
1	E	301	PRO
1	E	303	SER
1	F	145	GLN
1	F	220	SER
1	F	221	THR
1	F	223	THR
1	F	229	SER
1	F	250	GLU
1	F	276	LEU
1	F	278	SER
1	F	301	PRO
1	F	303	SER
1	F	304	ASN
1	F	306	ALA
1	G	150	THR
1	G	187	GLN
1	G	221	THR
1	G	243	ASN
1	G	250	GLU
1	G	299	GLU
1	G	303	SER
1	G	304	ASN
1	G	321	ASP
1	A	132	ASN
1	A	243	ASN
1	A	275	MET
1	A	301	PRO
1	A	305	ILE
1	A	321	ASP
1	B	150	THR
1	B	243	ASN
1	B	321	ASP
1	D	186	THR
1	D	218	LYS
1	D	220	SER
1	D	248	ASP
1	D	249	SER
1	D	278	SER
1	D	303	SER
1	D	307	THR
1	D	321	ASP
1	E	132	ASN

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Mol	Chain	Res	Type
1	E	148	GLU
1	E	223	THR
1	E	228	ALA
1	E	243	ASN
1	E	304	ASN
1	E	321	ASP
1	G	218	LYS
1	G	220	SER
1	G	248	ASP
1	G	308	LEU
1	A	219	SER
1	A	249	SER
1	A	262	SER
1	A	264	ASP
1	A	276	LEU
1	B	142	PRO
1	B	148	GLU
1	B	182	ASN
1	B	229	SER
1	B	232	ALA
1	B	303	SER
1	D	250	GLU
1	E	246	THR
1	E	281	ILE
1	F	142	PRO
1	F	150	THR
1	F	205	LEU
1	F	243	ASN
1	G	145	GLN
1	G	229	SER
1	G	262	SER
1	G	306	ALA
1	A	263	TYR
1	A	299	GLU
1	A	308	LEU
1	B	139	ASN
1	B	140	PRO
1	B	200	SER
1	B	227	VAL
1	B	248	ASP
1	D	187	GLN
1	D	229	SER

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Mol	Chain	Res	Type
1	D	263	TYR
1	E	229	SER
1	E	262	SER
1	E	276	LEU
1	E	308	LEU
1	F	140	PRO
1	F	227	VAL
1	G	249	SER
1	G	307	THR
1	A	229	SER
1	B	249	SER
1	B	299	GLU
1	E	249	SER
1	F	139	ASN
1	F	321	ASP
1	G	276	LEU
1	F	148	GLU
1	F	204	LEU
1	G	263	TYR
1	A	142	PRO
1	E	305	ILE
1	F	305	ILE
1	G	139	ASN
1	G	305	ILE
1	B	146	ILE
1	G	140	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	175/196 (89%)	155 (89%)	20 (11%)	5	14
1	B	175/196 (89%)	160 (91%)	15 (9%)	10	25
1	D	175/196 (89%)	153 (87%)	22 (13%)	4	11
1	E	175/196 (89%)	155 (89%)	20 (11%)	5	14

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	175/196 (89%)	153 (87%)	22 (13%)	4	11
1	G	175/196 (89%)	162 (93%)	13 (7%)	13	32
All	All	1050/1176 (89%)	938 (89%)	112 (11%)	6	16

All (112) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	143	ASN
1	A	145	GLN
1	A	163	LYS
1	A	172	VAL
1	A	174	LEU
1	A	188	LYS
1	A	217	ASN
1	A	221	THR
1	A	251	ASN
1	A	262	SER
1	A	263	TYR
1	A	265	ARG
1	A	270	LEU
1	A	272	ILE
1	A	275	MET
1	A	280	MET
1	A	284	ASN
1	A	298	SER
1	A	302	GLU
1	A	312	PRO
1	B	143	ASN
1	B	145	GLN
1	B	150	THR
1	B	161	LEU
1	B	172	VAL
1	B	174	LEU
1	B	188	LYS
1	B	221	THR
1	B	251	ASN
1	B	265	ARG
1	B	270	LEU
1	B	275	MET
1	B	284	ASN
1	B	301	PRO

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Mol	Chain	Res	Type
1	B	302	GLU
1	D	140	PRO
1	D	143	ASN
1	D	148	GLU
1	D	161	LEU
1	D	167	LEU
1	D	174	LEU
1	D	178	SER
1	D	180	THR
1	D	189	THR
1	D	200	SER
1	D	209	SER
1	D	213	ILE
1	D	217	ASN
1	D	251	ASN
1	D	261	THR
1	D	262	SER
1	D	265	ARG
1	D	270	LEU
1	D	275	MET
1	D	284	ASN
1	D	290	GLN
1	D	307	THR
1	E	143	ASN
1	E	144	CYS
1	E	145	GLN
1	E	172	VAL
1	E	173	SER
1	E	174	LEU
1	E	188	LYS
1	E	217	ASN
1	E	221	THR
1	E	251	ASN
1	E	257	CYS
1	E	263	TYR
1	E	265	ARG
1	E	270	LEU
1	E	272	ILE
1	E	275	MET
1	E	284	ASN
1	E	301	PRO
1	E	302	GLU

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Mol	Chain	Res	Type
1	E	307	THR
1	F	143	ASN
1	F	145	GLN
1	F	150	THR
1	F	161	LEU
1	F	172	VAL
1	F	174	LEU
1	F	179	ASP
1	F	188	LYS
1	F	205	LEU
1	F	217	ASN
1	F	221	THR
1	F	227	VAL
1	F	251	ASN
1	F	265	ARG
1	F	270	LEU
1	F	275	MET
1	F	280	MET
1	F	284	ASN
1	F	301	PRO
1	F	302	GLU
1	F	307	THR
1	F	319	THR
1	G	140	PRO
1	G	143	ASN
1	G	148	GLU
1	G	174	LEU
1	G	200	SER
1	G	217	ASN
1	G	251	ASN
1	G	262	SER
1	G	265	ARG
1	G	270	LEU
1	G	272	ILE
1	G	275	MET
1	G	284	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (71) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	132	ASN
1	A	139	ASN

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Mol	Chain	Res	Type
1	A	143	ASN
1	A	145	GLN
1	A	153	ASN
1	A	169	ASN
1	A	187	GLN
1	A	191	ASN
1	A	203	ASN
1	A	217	ASN
1	A	251	ASN
1	A	284	ASN
1	A	290	GLN
1	A	296	ASN
1	B	132	ASN
1	B	143	ASN
1	B	145	GLN
1	B	164	ASN
1	B	187	GLN
1	B	193	GLN
1	B	251	ASN
1	B	284	ASN
1	B	290	GLN
1	D	132	ASN
1	D	143	ASN
1	D	145	GLN
1	D	153	ASN
1	D	164	ASN
1	D	182	ASN
1	D	203	ASN
1	D	217	ASN
1	D	243	ASN
1	D	251	ASN
1	D	284	ASN
1	D	290	GLN
1	D	296	ASN
1	E	132	ASN
1	E	139	ASN
1	E	143	ASN
1	E	145	GLN
1	E	164	ASN
1	E	187	GLN
1	E	191	ASN
1	E	203	ASN

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Mol	Chain	Res	Type
1	E	217	ASN
1	E	251	ASN
1	E	284	ASN
1	E	290	GLN
1	E	294	ASN
1	F	132	ASN
1	F	143	ASN
1	F	145	GLN
1	F	193	GLN
1	F	203	ASN
1	F	217	ASN
1	F	251	ASN
1	F	284	ASN
1	F	290	GLN
1	F	296	ASN
1	G	132	ASN
1	G	143	ASN
1	G	164	ASN
1	G	191	ASN
1	G	203	ASN
1	G	217	ASN
1	G	243	ASN
1	G	251	ASN
1	G	284	ASN
1	G	290	GLN
1	G	294	ASN
1	G	296	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	192/220 (87%)	-0.75	2 (1%) 79 78	33, 72, 100, 108	66 (34%)
1	B	192/220 (87%)	-0.74	0 100 100	38, 76, 100, 108	71 (36%)
1	D	192/220 (87%)	-0.71	2 (1%) 79 78	42, 82, 103, 108	62 (32%)
1	E	192/220 (87%)	-0.73	2 (1%) 79 78	42, 72, 99, 106	68 (35%)
1	F	192/220 (87%)	-0.76	3 (1%) 70 68	38, 73, 102, 109	63 (32%)
1	G	192/220 (87%)	-0.64	1 (0%) 87 86	50, 81, 102, 108	63 (32%)
All	All	1152/1320 (87%)	-0.72	10 (0%) 81 80	33, 76, 102, 109	393 (34%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	246	THR	5.5
1	F	227	VAL	3.6
1	G	221	THR	3.1
1	D	223	THR	2.9
1	A	246	THR	2.7
1	D	238	THR	2.6
1	E	246	THR	2.5
1	F	223	THR	2.2
1	E	305	ILE	2.2
1	A	225	GLU	2.1

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

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.