



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 25, 2026 – 10:16 AM UTC

PDB ID : 6KQR / pdb_00006kqr
Title : A pre-assembled molecular-helical Cascade backbone of Csy3 subunits from *Zymomonas mobilis*
Authors : Gu, D.H.; Ha, S.C.; Kim, J.S.
Deposited on : 2019-08-18
Resolution : 2.90 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

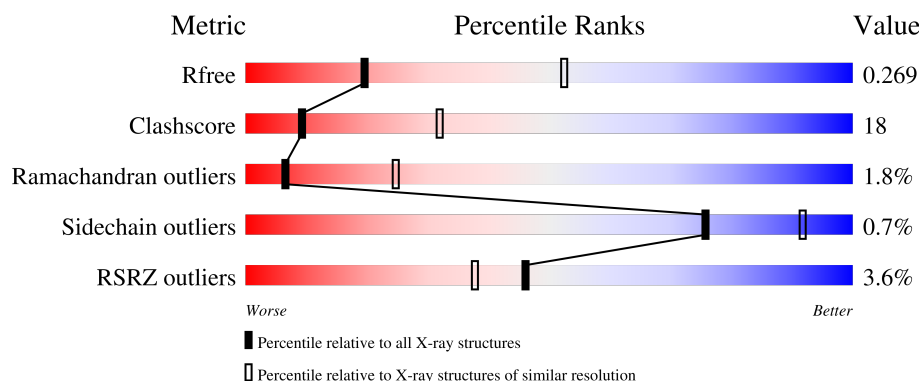
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	348	
1	B	348	
1	C	348	
1	D	348	
1	E	348	

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Mol	Chain	Length	Quality of chain
1	F	348	3% 64% 30%
1	G	348	4% 67% 28%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 18135 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 CRISPR-associated protein Csy3 family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	333	Total	C	N	O	S	0	0	0
			2589	1643	452	491	3			
1	B	333	Total	C	N	O	S	0	0	0
			2589	1643	452	491	3			
1	C	333	Total	C	N	O	S	0	0	0
			2589	1643	452	491	3			
1	D	333	Total	C	N	O	S	0	0	0
			2589	1643	452	491	3			
1	E	334	Total	C	N	O	S	0	0	0
			2593	1645	453	492	3			
1	F	334	Total	C	N	O	S	0	0	0
			2593	1645	453	492	3			
1	G	334	Total	C	N	O	S	0	0	0
			2593	1645	453	492	3			

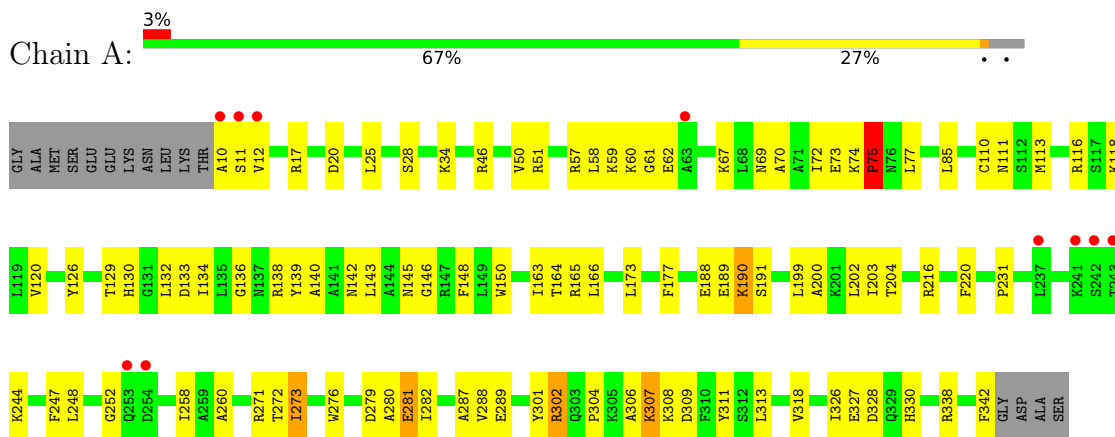
There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP A0A2R4RXD4
A	0	ALA	-	expression tag	UNP A0A2R4RXD4
B	-1	GLY	-	expression tag	UNP A0A2R4RXD4
B	0	ALA	-	expression tag	UNP A0A2R4RXD4
C	-1	GLY	-	expression tag	UNP A0A2R4RXD4
C	0	ALA	-	expression tag	UNP A0A2R4RXD4
D	-1	GLY	-	expression tag	UNP A0A2R4RXD4
D	0	ALA	-	expression tag	UNP A0A2R4RXD4
E	-1	GLY	-	expression tag	UNP A0A2R4RXD4
E	0	ALA	-	expression tag	UNP A0A2R4RXD4
F	-1	GLY	-	expression tag	UNP A0A2R4RXD4
F	0	ALA	-	expression tag	UNP A0A2R4RXD4
G	-1	GLY	-	expression tag	UNP A0A2R4RXD4
G	0	ALA	-	expression tag	UNP A0A2R4RXD4

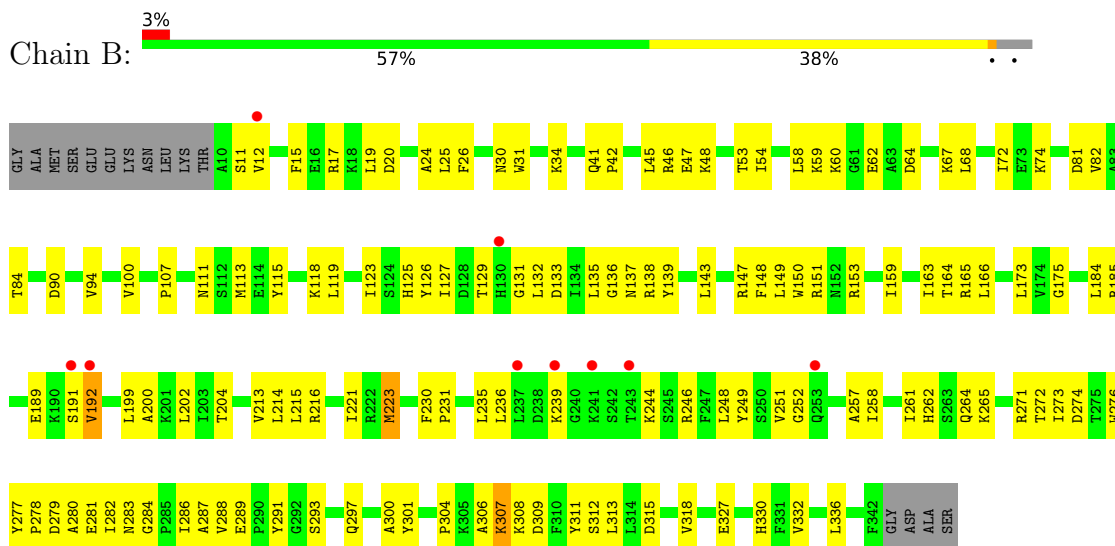
3 Residue-property plots

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.

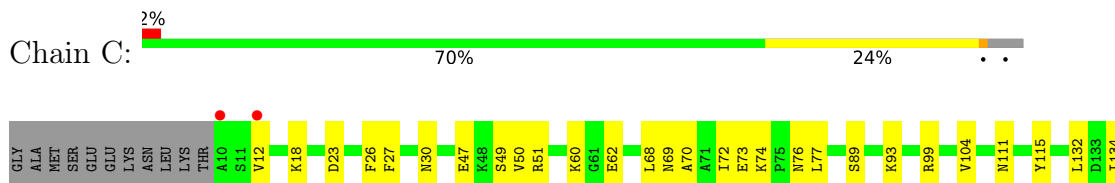
- Molecule 1: CRISPR-associated protein Csy3 family

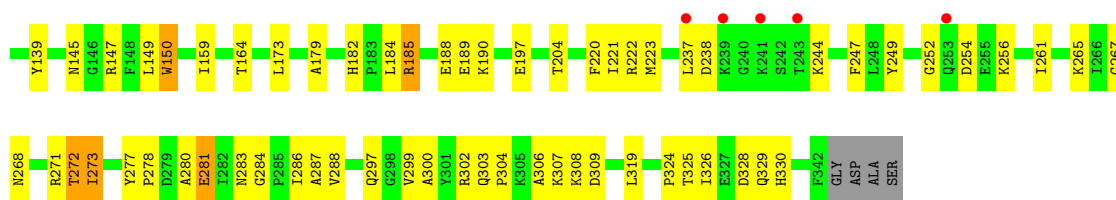


- Molecule 1: CRISPR-associated protein Csy3 family

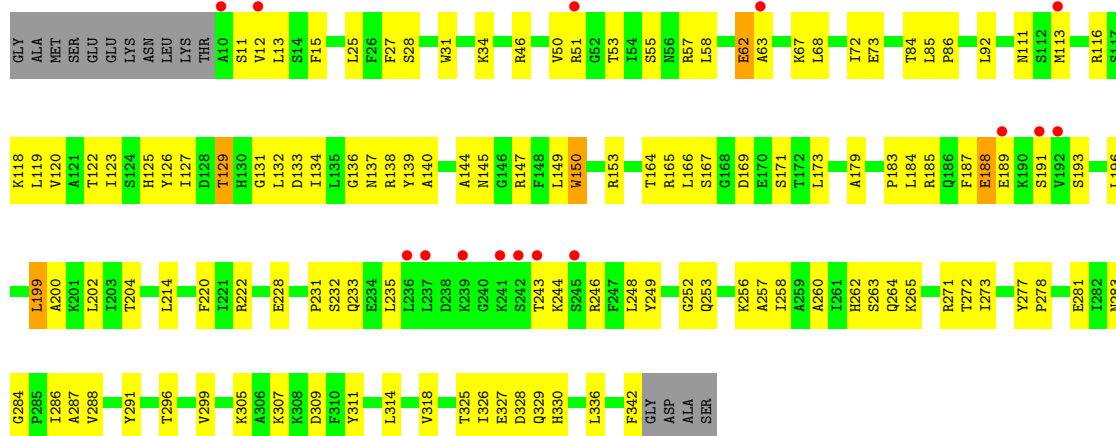


- Molecule 1: CRISPR-associated protein Csy3 family

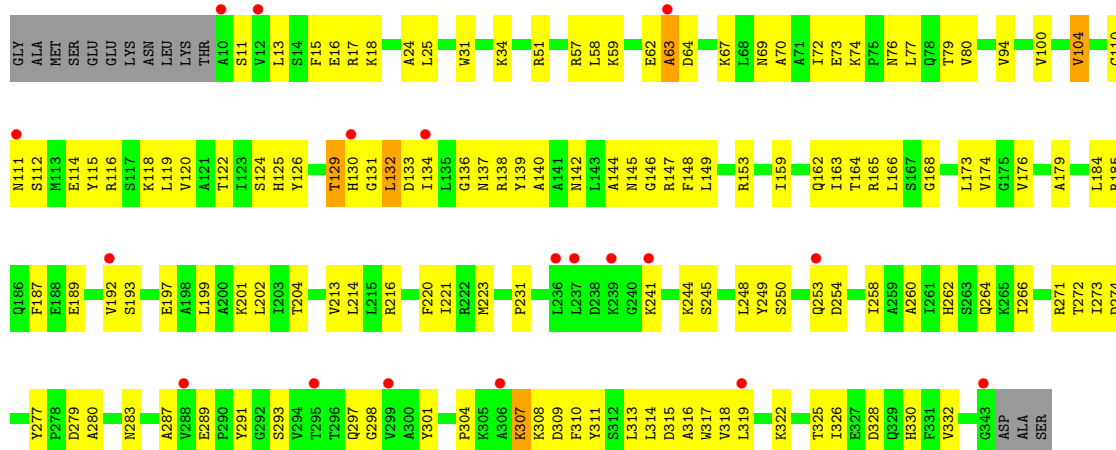




• Molecule 1: CRISPR-associated protein Csy3 family



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• Molecule 1: CRISPR-associated protein Csy3 family





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	82.18Å 116.16Å 115.94Å 99.37° 103.12° 103.35°	Depositor
Resolution (Å)	15.00 – 2.90 15.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.2 (15.00-2.90) 97.4 (15.00-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.214 , 0.267 0.215 , 0.269	Depositor DCC
R_{free} test set	3417 reflections (3.89%)	wwPDB-VP
Wilson B-factor (Å ²)	67.3	Xtriage
Anisotropy	0.783	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 54.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.012 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	18135	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.79% 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	0.56	1/2640 (0.0%)	0.73	1/3575 (0.0%)
1	B	0.41	0/2640	0.66	0/3575
1	C	0.49	0/2640	0.69	0/3575
1	D	0.64	1/2640 (0.0%)	0.77	3/3575 (0.1%)
1	E	0.45	0/2644	0.71	1/3580 (0.0%)
1	F	0.47	0/2644	0.72	0/3580
1	G	0.47	1/2644 (0.0%)	0.68	0/3580
All	All	0.51	3/18492 (0.0%)	0.71	5/25040 (0.0%)

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	C	0	1
1	D	0	1
All	All	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	86	PRO	N-CD	-22.30	1.16	1.47
1	A	75	PRO	N-CD	-16.33	1.24	1.47
1	G	75	PRO	N-CD	-9.78	1.34	1.47

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	189	GLU	N-CA-C	8.97	129.91	110.80
1	D	53	THR	CA-C-N	-6.19	113.13	121.80
1	D	53	THR	C-N-CA	-6.19	113.13	121.80
1	A	75	PRO	CA-N-CD	5.42	119.59	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	254	ASP	CB-CA-C	-5.34	109.96	117.23

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	272	THR	Peptide
1	D	188	GLU	Peptide

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	2589	0	2591	96	1
1	B	2589	0	2591	127	0
1	C	2589	0	2591	76	0
1	D	2589	0	2591	99	1
1	E	2593	0	2594	131	0
1	F	2593	0	2594	110	0
1	G	2593	0	2594	97	0
All	All	18135	0	18146	656	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 18.

All (656) 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:G:74:LYS:CD	1:G:75:PRO:HD2	1.62	1.30
1:G:74:LYS:HD2	1:G:75:PRO:CD	1.63	1.27
1:A:74:LYS:CD	1:A:75:PRO:HD2	1.79	1.13
1:B:47:GLU:OE2	1:B:244:LYS:NZ	1.82	1.11
1:A:74:LYS:HD2	1:A:75:PRO:HD2	1.19	1.09
1:B:123:ILE:O	1:B:127:ILE:HD12	1.53	1.07
1:E:51:ARG:HH21	1:F:265:LYS:HD3	1.14	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:11:SER:HB3	1:E:112:SER:HB3	1.38	1.04
1:F:77:LEU:HD11	1:G:235:LEU:HD12	1.39	1.03
1:G:311:TYR:HH	1:G:342:PHE:HD2	1.04	1.02
1:G:60:LYS:HG3	1:G:61:GLY:H	1.19	1.02
1:F:134:ILE:HG21	1:F:330:HIS:CD2	1.96	0.99
1:E:62:GLU:HA	1:E:67:LYS:HD2	1.45	0.95
1:F:51:ARG:HH22	1:G:265:LYS:HB2	1.30	0.95
1:A:311:TYR:HH	1:A:342:PHE:HD1	0.99	0.94
1:G:51:ARG:H	1:G:51:ARG:HD3	1.35	0.91
1:F:77:LEU:CD1	1:G:235:LEU:HD12	2.02	0.90
1:C:47:GLU:OE2	1:C:244:LYS:NZ	2.06	0.88
1:A:74:LYS:HD2	1:A:75:PRO:CD	2.03	0.88
1:C:51:ARG:HD3	1:C:51:ARG:H	1.40	0.87
1:B:279:ASP:HB3	1:B:282:ILE:HD11	1.58	0.85
1:A:136:GLY:HA3	1:A:204:THR:HG22	1.60	0.83
1:G:136:GLY:HA3	1:G:204:THR:HG23	1.58	0.83
1:G:310:PHE:HD2	1:G:311:TYR:CE1	1.97	0.83
1:A:134:ILE:HD12	1:A:134:ILE:H	1.42	0.83
1:A:59:LYS:H	1:A:59:LYS:HD3	1.45	0.82
1:E:272:THR:HA	1:E:287:ALA:HA	1.61	0.82
1:G:13:LEU:HD23	1:G:15:PHE:HE1	1.44	0.81
1:G:13:LEU:HD23	1:G:15:PHE:CE1	2.16	0.81
1:F:271:ARG:NH1	1:F:289:GLU:OE2	2.14	0.81
1:D:165:ARG:NH1	1:D:167:SER:OG	2.13	0.81
1:F:65:PRO:HA	1:G:303:GLN:NE2	1.96	0.81
1:G:60:LYS:HG3	1:G:61:GLY:N	1.96	0.80
1:D:185:ARG:H	1:D:185:ARG:HD3	1.46	0.80
1:G:310:PHE:HD2	1:G:311:TYR:CD1	2.01	0.79
1:A:62:GLU:HG3	1:A:67:LYS:HG2	1.64	0.79
1:B:12:VAL:HB	1:B:111:ASN:HB2	1.64	0.79
1:E:136:GLY:HA3	1:E:204:THR:HG23	1.64	0.79
1:F:65:PRO:HA	1:G:303:GLN:HE22	1.48	0.79
1:A:163:ILE:HD12	1:A:199:LEU:HD11	1.65	0.79
1:F:134:ILE:HG21	1:F:330:HIS:HD2	1.46	0.78
1:E:51:ARG:HH22	1:F:265:LYS:HB2	1.49	0.78
1:C:280:ALA:O	1:C:281:GLU:HG2	1.84	0.78
1:A:110:CYS:SG	1:A:113:MET:CE	2.72	0.77
1:A:326:ILE:HD12	1:A:326:ILE:H	1.49	0.77
1:B:136:GLY:HA3	1:B:204:THR:HG23	1.65	0.77
1:A:77:LEU:HD11	1:B:235:LEU:HD12	1.66	0.76
1:B:123:ILE:HG22	1:B:127:ILE:HD11	1.67	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:51:ARG:NH2	1:F:265:LYS:HD3	1.95	0.76
1:D:228:GLU:OE1	1:E:17:ARG:NH2	2.18	0.76
1:C:267:GLY:C	1:C:271:ARG:HH21	1.94	0.76
1:A:51:ARG:H	1:A:51:ARG:HD3	1.51	0.75
1:C:268:ASN:N	1:C:271:ARG:NH2	2.35	0.74
1:C:190:LYS:HG2	1:C:197:GLU:HG3	1.68	0.73
1:A:51:ARG:HH12	1:B:262:HIS:CD2	2.07	0.73
1:B:133:ASP:HA	1:B:204:THR:HG21	1.70	0.73
1:B:279:ASP:O	1:B:281:GLU:N	2.22	0.73
1:E:51:ARG:NH2	1:F:265:LYS:HB2	2.03	0.73
1:A:72:ILE:HD13	1:B:293:SER:HB2	1.71	0.72
1:B:306:ALA:O	1:B:308:LYS:N	2.22	0.72
1:E:145:ASN:O	1:E:153:ARG:NH1	2.22	0.72
1:F:57:ARG:HG2	1:G:311:TYR:CD2	2.25	0.72
1:E:100:VAL:HB	1:E:213:VAL:HG11	1.71	0.71
1:F:46:ARG:NH2	1:G:249:TYR:OH	2.22	0.71
1:E:139:TYR:CD1	1:E:273:ILE:HD11	2.25	0.71
1:F:338:ARG:HD3	1:F:339:GLY:O	1.91	0.71
1:C:267:GLY:C	1:C:271:ARG:NH2	2.50	0.70
1:G:11:SER:HB2	1:G:13:LEU:HD13	1.74	0.70
1:A:70:ALA:O	1:A:73:GLU:OE2	2.10	0.70
1:E:79:THR:HB	1:E:244:LYS:HD2	1.72	0.70
1:C:51:ARG:HH21	1:D:265:LYS:HD3	1.56	0.70
1:C:51:ARG:HH22	1:D:265:LYS:HB2	1.57	0.70
1:A:51:ARG:HH12	1:B:262:HIS:HD2	1.40	0.69
1:A:74:LYS:HD3	1:A:75:PRO:HD2	1.71	0.69
1:C:134:ILE:HD12	1:C:134:ILE:H	1.58	0.69
1:E:72:ILE:HD12	1:F:300:ALA:HB2	1.74	0.69
1:D:272:THR:HA	1:D:287:ALA:HA	1.73	0.69
1:G:42:PRO:HB3	1:G:258:ILE:HD13	1.74	0.69
1:A:51:ARG:HH21	1:B:265:LYS:HD3	1.58	0.69
1:B:42:PRO:HB2	1:B:258:ILE:CD1	2.23	0.69
1:A:110:CYS:SG	1:A:113:MET:HE1	2.32	0.69
1:B:42:PRO:HB2	1:B:258:ILE:HD13	1.74	0.69
1:B:42:PRO:CB	1:B:258:ILE:CD1	2.72	0.68
1:D:55:SER:O	1:E:311:TYR:OH	2.04	0.68
1:B:153:ARG:HH12	1:B:184:LEU:HD13	1.59	0.68
1:D:140:ALA:HB1	1:D:199:LEU:HD13	1.74	0.68
1:E:72:ILE:CD1	1:F:300:ALA:HB2	2.24	0.68
1:G:133:ASP:HA	1:G:204:THR:HG21	1.74	0.68
1:B:166:LEU:HD11	1:B:216:ARG:HB2	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:133:ASP:HA	1:D:204:THR:HG21	1.75	0.67
1:A:166:LEU:HD21	1:A:216:ARG:HH21	1.60	0.67
1:D:12:VAL:HB	1:D:111:ASN:HB2	1.74	0.67
1:C:164:THR:HG22	1:C:173:LEU:HA	1.77	0.67
1:B:185:ARG:H	1:B:185:ARG:HD3	1.60	0.67
1:E:187:PHE:O	1:E:189:GLU:OE2	2.12	0.67
1:A:166:LEU:HD21	1:A:216:ARG:NH2	2.11	0.66
1:D:136:GLY:HA3	1:D:204:THR:HG22	1.78	0.66
1:G:119:LEU:HD12	1:G:318:VAL:HG21	1.78	0.66
1:A:51:ARG:HH22	1:B:265:LYS:HB2	1.61	0.66
1:D:134:ILE:HG21	1:D:330:HIS:CE1	2.31	0.66
1:G:310:PHE:CD2	1:G:311:TYR:CD1	2.84	0.66
1:E:125:HIS:O	1:E:129:THR:HG22	1.95	0.66
1:F:51:ARG:H	1:F:51:ARG:HD3	1.58	0.66
1:A:51:ARG:NH2	1:B:265:LYS:HB2	2.10	0.66
1:F:51:ARG:NH2	1:G:265:LYS:HD3	2.11	0.66
1:F:237:LEU:HD21	1:F:245:SER:O	1.97	0.65
1:C:278:PRO:HD3	1:C:308:LYS:HE2	1.78	0.65
1:B:125:HIS:O	1:B:129:THR:HG23	1.97	0.65
1:E:69:ASN:ND2	1:F:299:VAL:HA	2.12	0.65
1:A:133:ASP:HA	1:A:204:THR:HG21	1.77	0.65
1:C:221:ILE:HG22	1:C:223:MET:HG2	1.78	0.65
1:E:140:ALA:HB1	1:E:199:LEU:HD23	1.79	0.65
1:G:13:LEU:HD21	1:G:119:LEU:HD22	1.78	0.65
1:F:278:PRO:HD3	1:F:308:LYS:HE2	1.77	0.65
1:A:164:THR:HG22	1:A:173:LEU:HA	1.78	0.65
1:C:68:LEU:O	1:C:72:ILE:HD12	1.98	0.64
1:E:18:LYS:HE3	1:E:104:VAL:HA	1.80	0.64
1:A:311:TYR:HE1	1:A:342:PHE:HB3	1.63	0.64
1:B:165:ARG:HE	1:B:202:LEU:HD11	1.62	0.64
1:D:126:TYR:HD2	1:D:127:ILE:HD12	1.63	0.64
1:C:244:LYS:HB2	1:C:247:PHE:CZ	2.32	0.64
1:A:140:ALA:HB2	1:A:203:ILE:HD12	1.78	0.64
1:C:51:ARG:NH2	1:D:265:LYS:HB2	2.13	0.64
1:G:185:ARG:H	1:G:185:ARG:HD3	1.62	0.64
1:E:72:ILE:HD13	1:F:293:SER:HB2	1.79	0.64
1:C:73:GLU:HB3	1:E:111:ASN:HD21	1.63	0.63
1:D:153:ARG:NH1	1:D:184:LEU:HD11	2.13	0.63
1:E:316:ALA:HB1	1:E:322:LYS:HB2	1.81	0.63
1:F:165:ARG:HG3	1:F:202:LEU:HD21	1.79	0.63
1:A:189:GLU:O	1:A:191:SER:N	2.31	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:60:LYS:O	1:C:62:GLU:HA	1.99	0.63
1:D:232:SER:HB3	1:D:262:HIS:CD2	2.34	0.63
1:E:59:LYS:HD2	1:E:62:GLU:HG3	1.80	0.63
1:E:133:ASP:HA	1:E:204:THR:HG21	1.80	0.62
1:F:129:THR:OG1	1:F:130:HIS:N	2.31	0.62
1:D:123:ILE:HG22	1:D:127:ILE:HD13	1.81	0.62
1:D:311:TYR:HE1	1:D:342:PHE:HB3	1.64	0.62
1:D:200:ALA:O	1:D:204:THR:HG23	1.99	0.62
1:A:34:LYS:HG2	1:A:220:PHE:CD2	2.34	0.62
1:E:120:VAL:O	1:E:124:SER:OG	2.15	0.62
1:F:51:ARG:NH2	1:G:265:LYS:HB2	2.09	0.62
1:B:15:PHE:CE2	1:B:336:LEU:HD22	2.35	0.62
1:G:164:THR:HG22	1:G:173:LEU:HA	1.81	0.62
1:G:234:GLU:C	1:G:235:LEU:HD23	2.25	0.62
1:B:60:LYS:H	1:B:60:LYS:HD2	1.65	0.62
1:C:139:TYR:CE1	1:C:273:ILE:HD12	2.35	0.62
1:B:276:TRP:HB3	1:B:327:GLU:HB3	1.82	0.62
1:F:51:ARG:NH1	1:G:262:HIS:HD2	1.98	0.62
1:E:262:HIS:HB3	1:E:264:GLN:OE1	1.99	0.62
1:A:51:ARG:H	1:A:51:ARG:CD	2.12	0.61
1:B:15:PHE:CE1	1:B:107:PRO:HB3	2.35	0.61
1:D:147:ARG:NH2	1:D:184:LEU:O	2.33	0.61
1:D:11:SER:HB2	1:D:13:LEU:HD13	1.81	0.61
1:E:51:ARG:H	1:E:51:ARG:HD3	1.65	0.61
1:F:134:ILE:CG2	1:F:330:HIS:HD2	2.12	0.61
1:A:57:ARG:HG2	1:B:311:TYR:CE2	2.36	0.61
1:C:74:LYS:HD3	1:C:76:ASN:H	1.65	0.61
1:F:110:CYS:SG	1:F:113:MET:HE1	2.40	0.61
1:C:272:THR:HA	1:C:287:ALA:HA	1.81	0.61
1:D:139:TYR:CE1	1:D:273:ILE:HD12	2.35	0.61
1:A:189:GLU:H	1:A:189:GLU:CD	2.08	0.61
1:A:200:ALA:O	1:A:204:THR:HG23	2.01	0.61
1:C:252:GLY:O	1:C:256:LYS:HG2	2.00	0.61
1:F:51:ARG:HH21	1:G:265:LYS:HD3	1.63	0.61
1:D:164:THR:HG22	1:D:173:LEU:HA	1.83	0.60
1:G:54:ILE:HG13	1:G:72:ILE:HD13	1.83	0.60
1:G:277:TYR:OH	1:G:283:ASN:HB3	2.00	0.60
1:C:12:VAL:H	1:C:111:ASN:CB	2.14	0.60
1:F:51:ARG:HH12	1:G:262:HIS:HD2	1.48	0.60
1:B:283:ASN:OD1	1:B:301:TYR:OH	2.20	0.60
1:C:18:LYS:HE3	1:C:104:VAL:HA	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:221:ILE:HG22	1:E:223:MET:HG2	1.83	0.60
1:F:164:THR:HG22	1:F:173:LEU:HA	1.84	0.60
1:G:51:ARG:H	1:G:51:ARG:CD	2.12	0.60
1:C:222:ARG:NH2	1:D:169:ASP:OD1	2.34	0.59
1:C:12:VAL:H	1:C:111:ASN:HB2	1.68	0.59
1:D:138:ARG:HB3	1:D:273:ILE:HG13	1.83	0.59
1:F:272:THR:HA	1:F:287:ALA:HA	1.83	0.59
1:C:134:ILE:HD13	1:C:330:HIS:HE1	1.66	0.59
1:C:324:PRO:HG2	1:C:329:GLN:HG2	1.84	0.59
1:D:51:ARG:H	1:D:51:ARG:HD3	1.67	0.59
1:B:272:THR:HA	1:B:287:ALA:HA	1.84	0.59
1:C:325:THR:HG23	1:C:328:ASP:H	1.68	0.59
1:E:119:LEU:HD12	1:E:318:VAL:HG21	1.85	0.59
1:E:129:THR:HG23	1:E:130:HIS:H	1.67	0.59
1:E:166:LEU:HB2	1:E:214:LEU:HD23	1.85	0.58
1:B:42:PRO:HB3	1:B:258:ILE:CD1	2.33	0.58
1:B:279:ASP:CB	1:B:282:ILE:HD11	2.29	0.58
1:F:149:LEU:HB2	1:F:153:ARG:HG3	1.86	0.58
1:G:244:LYS:HD2	1:G:247:PHE:CE2	2.38	0.58
1:C:297:GLN:HB2	1:C:299:VAL:HG12	1.84	0.58
1:F:27:PHE:HB2	1:F:93:LYS:HB3	1.85	0.58
1:G:11:SER:HB2	1:G:13:LEU:CD1	2.32	0.58
1:A:61:GLY:HA2	1:A:62:GLU:HB3	1.85	0.58
1:G:189:GLU:OE2	1:G:189:GLU:N	2.28	0.58
1:C:306:ALA:O	1:C:308:LYS:N	2.36	0.57
1:E:271:ARG:NH1	1:E:289:GLU:OE2	2.37	0.57
1:F:51:ARG:HH22	1:G:265:LYS:CB	2.12	0.57
1:D:116:ARG:O	1:D:120:VAL:HG23	2.04	0.57
1:E:142:ASN:OD1	1:E:145:ASN:ND2	2.38	0.57
1:G:182:HIS:ND1	1:G:188:GLU:OE1	2.31	0.57
1:G:310:PHE:CD2	1:G:311:TYR:CE1	2.88	0.57
1:D:119:LEU:HD11	1:D:314:LEU:HD11	1.87	0.57
1:F:57:ARG:HG2	1:G:311:TYR:HD2	1.70	0.57
1:A:306:ALA:O	1:A:308:LYS:N	2.37	0.57
1:C:182:HIS:ND1	1:C:188:GLU:OE2	2.37	0.57
1:A:46:ARG:NH2	1:B:249:TYR:OH	2.38	0.56
1:A:126:TYR:OH	1:A:330:HIS:ND1	2.32	0.56
1:B:191:SER:O	1:B:192:VAL:HG22	2.04	0.56
1:B:236:LEU:HD12	1:B:239:LYS:CB	2.35	0.56
1:A:306:ALA:O	1:A:308:LYS:HG2	2.05	0.56
1:E:69:ASN:HD22	1:F:299:VAL:HA	1.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:131:GLY:O	1:E:133:ASP:N	2.36	0.56
1:F:100:VAL:HB	1:F:213:VAL:CG2	2.35	0.56
1:A:51:ARG:NH1	1:B:262:HIS:HD2	2.03	0.56
1:B:123:ILE:HG22	1:B:127:ILE:CD1	2.35	0.56
1:E:51:ARG:H	1:E:51:ARG:CD	2.18	0.56
1:G:153:ARG:CZ	1:G:184:LEU:HD21	2.36	0.56
1:D:25:LEU:HD13	1:D:258:ILE:HB	1.88	0.56
1:F:134:ILE:CG2	1:F:330:HIS:CD2	2.81	0.56
1:G:233:GLN:HB3	1:G:235:LEU:HD21	1.86	0.56
1:A:59:LYS:HE3	1:A:61:GLY:O	2.06	0.56
1:D:271:ARG:HG3	1:D:288:VAL:H	1.70	0.56
1:A:116:ARG:O	1:A:120:VAL:HG23	2.06	0.56
1:B:153:ARG:HH12	1:B:184:LEU:CD1	2.19	0.56
1:E:57:ARG:HE	1:F:343:GLY:HA2	1.69	0.55
1:B:279:ASP:HB3	1:B:282:ILE:CD1	2.31	0.55
1:D:314:LEU:O	1:D:318:VAL:HG22	2.06	0.55
1:B:123:ILE:C	1:B:127:ILE:HD12	2.30	0.55
1:C:73:GLU:CB	1:E:111:ASN:HD21	2.18	0.55
1:G:116:ARG:O	1:G:120:VAL:HG12	2.05	0.55
1:A:129:THR:OG1	1:A:130:HIS:N	2.34	0.55
1:D:58:LEU:HD11	1:D:68:LEU:HD12	1.87	0.55
1:A:139:TYR:CE1	1:A:273:ILE:HD12	2.41	0.55
1:B:62:GLU:HA	1:B:67:LYS:CE	2.37	0.55
1:F:38:LYS:O	1:F:38:LYS:HD3	2.07	0.55
1:E:142:ASN:HA	1:E:145:ASN:ND2	2.21	0.55
1:F:231:PRO:HG2	1:F:248:LEU:HD13	1.88	0.55
1:B:118:LYS:HG3	1:B:318:VAL:HG13	1.89	0.55
1:A:272:THR:HA	1:A:287:ALA:HA	1.89	0.55
1:B:25:LEU:HD21	1:B:251:VAL:CG2	2.37	0.55
1:F:244:LYS:HE2	1:F:247:PHE:CE1	2.42	0.55
1:A:70:ALA:HA	1:A:73:GLU:OE2	2.07	0.55
1:C:69:ASN:OD1	1:D:299:VAL:HA	2.07	0.55
1:E:115:TYR:HE1	1:E:119:LEU:HD13	1.72	0.55
1:A:326:ILE:H	1:A:326:ILE:CD1	2.18	0.54
1:F:134:ILE:HA	1:F:137:ASN:HB2	1.90	0.54
1:D:51:ARG:HD3	1:D:51:ARG:N	2.22	0.54
1:E:77:LEU:O	1:E:77:LEU:HD12	2.08	0.54
1:A:70:ALA:C	1:A:73:GLU:OE2	2.50	0.54
1:F:59:LYS:HE2	1:F:59:LYS:H	1.71	0.54
1:F:68:LEU:HD22	1:F:72:ILE:HD11	1.89	0.54
1:F:136:GLY:HA3	1:F:204:THR:CG2	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:244:LYS:HE2	1:F:247:PHE:CD1	2.42	0.54
1:C:147:ARG:NH1	1:C:184:LEU:HG	2.23	0.54
1:G:165:ARG:NH1	1:G:202:LEU:HD11	2.22	0.54
1:C:139:TYR:CD1	1:C:273:ILE:HD11	2.42	0.54
1:F:12:VAL:H	1:F:111:ASN:HB3	1.71	0.54
1:C:115:TYR:CE2	1:C:319:LEU:HD21	2.43	0.54
1:D:185:ARG:H	1:D:185:ARG:CD	2.17	0.54
1:G:166:LEU:HD21	1:G:216:ARG:NH1	2.23	0.54
1:C:189:GLU:CD	1:C:189:GLU:H	2.15	0.54
1:D:183:PRO:HD2	1:D:188:GLU:OE2	2.08	0.54
1:G:17:ARG:HD2	1:G:20:ASP:OD1	2.08	0.54
1:C:326:ILE:H	1:C:326:ILE:HD12	1.73	0.53
1:D:63:ALA:O	1:E:304:PRO:HG2	2.08	0.53
1:D:145:ASN:O	1:D:179:ALA:HB1	2.08	0.53
1:F:100:VAL:CG1	1:F:207:LEU:HD22	2.38	0.53
1:G:132:LEU:HD12	1:G:132:LEU:O	2.08	0.53
1:D:273:ILE:HG22	1:D:288:VAL:CG2	2.38	0.53
1:B:41:GLN:OE1	1:B:42:PRO:HD2	2.08	0.53
1:C:30:ASN:OD1	1:C:89:SER:HA	2.08	0.53
1:C:300:ALA:HB3	1:C:303:GLN:NE2	2.23	0.53
1:A:142:ASN:HA	1:A:145:ASN:OD1	2.09	0.53
1:C:280:ALA:O	1:C:281:GLU:CG	2.53	0.53
1:G:110:CYS:HB3	1:G:116:ARG:NH1	2.22	0.53
1:A:190:LYS:HG2	1:A:190:LYS:O	2.08	0.53
1:B:25:LEU:HD21	1:B:251:VAL:HG22	1.91	0.53
1:B:304:PRO:O	1:B:307:LYS:HD3	2.09	0.53
1:E:31:TRP:O	1:E:34:LYS:NZ	2.41	0.53
1:E:112:SER:OG	1:E:114:GLU:HG2	2.09	0.53
1:E:163:ILE:HG21	1:E:199:LEU:HD11	1.91	0.53
1:G:68:LEU:O	1:G:72:ILE:HG12	2.08	0.53
1:A:58:LEU:HD13	1:B:291:TYR:CZ	2.44	0.53
1:F:153:ARG:CZ	1:F:184:LEU:HD11	2.38	0.53
1:F:271:ARG:HG3	1:F:288:VAL:HB	1.90	0.53
1:E:110:CYS:HB3	1:E:116:ARG:CZ	2.39	0.53
1:C:74:LYS:HD3	1:C:76:ASN:N	2.24	0.52
1:E:69:ASN:HD21	1:F:299:VAL:HG23	1.73	0.52
1:E:277:TYR:CZ	1:E:280:ALA:HA	2.44	0.52
1:B:131:GLY:C	1:B:133:ASP:H	2.18	0.52
1:E:326:ILE:O	1:E:330:HIS:ND1	2.42	0.52
1:G:143:LEU:HD12	1:G:199:LEU:HD21	1.90	0.52
1:B:47:GLU:CD	1:B:244:LYS:HZ1	2.14	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:325:THR:O	1:C:329:GLN:HG3	2.09	0.52
1:E:119:LEU:HD11	1:E:314:LEU:HD21	1.90	0.52
1:G:165:ARG:HD2	1:G:202:LEU:HD21	1.91	0.52
1:G:310:PHE:CD2	1:G:311:TYR:HD1	2.27	0.52
1:A:304:PRO:HA	1:A:309:ASP:OD1	2.08	0.52
1:D:134:ILE:HG21	1:D:330:HIS:ND1	2.24	0.52
1:E:310:PHE:CE1	1:E:332:VAL:HG23	2.45	0.52
1:B:231:PRO:HG2	1:B:248:LEU:HD13	1.90	0.52
1:D:171:SER:O	1:D:171:SER:OG	2.26	0.52
1:E:297:GLN:HA	1:F:110:CYS:O	2.09	0.52
1:A:136:GLY:HA3	1:A:204:THR:CG2	2.37	0.52
1:E:164:THR:HG22	1:E:173:LEU:HA	1.92	0.52
1:E:316:ALA:O	1:E:322:LYS:N	2.41	0.52
1:G:15:PHE:HZ	1:G:119:LEU:HD21	1.75	0.52
1:C:51:ARG:HD3	1:C:51:ARG:N	2.16	0.52
1:C:150:TRP:CE2	1:C:265:LYS:HE3	2.45	0.52
1:C:244:LYS:HD2	1:C:247:PHE:CE1	2.45	0.52
1:B:45:LEU:HD13	1:B:248:LEU:HB2	1.92	0.52
1:B:271:ARG:NH2	1:B:289:GLU:OE2	2.43	0.52
1:B:276:TRP:CH2	1:B:330:HIS:HB3	2.44	0.52
1:E:58:LEU:HB3	1:E:62:GLU:HB2	1.92	0.52
1:E:325:THR:HG23	1:E:328:ASP:HB2	1.92	0.52
1:G:262:HIS:HB3	1:G:264:GLN:OE1	2.09	0.52
1:A:57:ARG:NH1	1:B:315:ASP:OD2	2.43	0.51
1:B:153:ARG:NH1	1:B:184:LEU:CD1	2.73	0.51
1:B:163:ILE:HD12	1:B:199:LEU:HD11	1.92	0.51
1:E:139:TYR:CE1	1:E:273:ILE:CD1	2.93	0.51
1:B:31:TRP:CE2	1:B:34:LYS:HE3	2.45	0.51
1:B:202:LEU:HD23	1:B:202:LEU:O	2.11	0.51
1:E:174:VAL:HG11	1:E:199:LEU:HD12	1.93	0.51
1:A:326:ILE:HD12	1:A:326:ILE:N	2.23	0.51
1:E:62:GLU:O	1:E:64:ASP:N	2.43	0.51
1:F:262:HIS:HB3	1:F:264:GLN:OE1	2.11	0.51
1:B:58:LEU:HB3	1:B:62:GLU:HG3	1.93	0.51
1:B:118:LYS:HG3	1:B:318:VAL:CG1	2.41	0.51
1:E:116:ARG:O	1:E:120:VAL:HG23	2.10	0.51
1:F:50:VAL:HG22	1:G:264:GLN:NE2	2.26	0.51
1:G:80:VAL:HG12	1:G:246:ARG:HB3	1.92	0.51
1:G:150:TRP:CD2	1:G:265:LYS:HE3	2.45	0.51
1:B:47:GLU:CD	1:B:244:LYS:NZ	2.67	0.51
1:F:110:CYS:SG	1:F:116:ARG:NH1	2.83	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:64:ASP:HB3	1:G:67:LYS:HE3	1.92	0.51
1:A:50:VAL:HG22	1:B:264:GLN:NE2	2.27	0.50
1:D:139:TYR:CD1	1:D:273:ILE:HD12	2.46	0.50
1:B:53:THR:HG22	1:B:54:ILE:H	1.76	0.50
1:C:271:ARG:HG3	1:C:288:VAL:H	1.75	0.50
1:B:46:ARG:NH1	1:B:84:THR:OG1	2.44	0.50
1:G:223:MET:O	1:G:227:GLN:HG3	2.11	0.50
1:C:277:TYR:OH	1:C:284:GLY:O	2.29	0.50
1:E:15:PHE:HZ	1:E:119:LEU:HD21	1.76	0.50
1:E:31:TRP:NE1	1:E:34:LYS:HZ1	2.09	0.50
1:E:301:TYR:CD1	1:E:301:TYR:N	2.78	0.50
1:G:150:TRP:CE2	1:G:265:LYS:HE3	2.46	0.50
1:F:25:LEU:HD13	1:F:258:ILE:HB	1.92	0.50
1:C:139:TYR:CD1	1:C:273:ILE:CD1	2.95	0.50
1:A:59:LYS:CE	1:A:61:GLY:O	2.59	0.50
1:B:309:ASP:OD2	1:B:312:SER:OG	2.24	0.50
1:E:148:PHE:HB2	1:E:266:ILE:HD13	1.93	0.50
1:A:271:ARG:NH1	1:A:289:GLU:OE2	2.45	0.49
1:F:72:ILE:CD1	1:G:300:ALA:HB2	2.41	0.49
1:A:313:LEU:HD13	1:A:328:ASP:HB3	1.94	0.49
1:B:64:ASP:OD2	1:B:67:LYS:HD3	2.12	0.49
1:B:185:ARG:HD3	1:B:185:ARG:N	2.26	0.49
1:F:277:TYR:OH	1:F:284:GLY:O	2.18	0.49
1:A:25:LEU:HD22	1:A:260:ALA:HB2	1.95	0.49
1:B:42:PRO:HB3	1:B:258:ILE:HD12	1.95	0.49
1:A:17:ARG:NE	1:A:20:ASP:OD1	2.46	0.49
1:E:100:VAL:HB	1:E:213:VAL:CG1	2.40	0.49
1:E:145:ASN:HB2	1:E:147:ARG:H	1.78	0.49
1:D:325:THR:OG1	1:D:327:GLU:OE1	2.31	0.49
1:F:133:ASP:HA	1:F:204:THR:HG21	1.94	0.49
1:G:85:LEU:HD21	1:G:92:LEU:HB2	1.94	0.49
1:G:306:ALA:HB1	1:G:308:LYS:CE	2.43	0.49
1:B:143:LEU:HD13	1:B:163:ILE:HD13	1.95	0.49
1:E:58:LEU:HG	1:F:291:TYR:CZ	2.48	0.49
1:B:147:ARG:NH2	1:B:184:LEU:O	2.46	0.49
1:G:142:ASN:HA	1:G:145:ASN:OD1	2.13	0.49
1:F:311:TYR:HE1	1:F:342:PHE:HB3	1.77	0.49
1:C:132:LEU:O	1:C:204:THR:HG23	2.12	0.49
1:F:57:ARG:HG2	1:G:311:TYR:CE2	2.48	0.49
1:F:131:GLY:O	1:F:133:ASP:N	2.43	0.49
1:F:136:GLY:HA3	1:F:204:THR:HG22	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:SER:HB2	1:A:85:LEU:HD12	1.95	0.48
1:B:68:LEU:O	1:B:72:ILE:HG12	2.12	0.48
1:B:332:VAL:O	1:B:336:LEU:HG	2.13	0.48
1:G:310:PHE:HD2	1:G:311:TYR:HE1	1.56	0.48
1:E:231:PRO:HG2	1:E:248:LEU:HD13	1.95	0.48
1:E:184:LEU:HB2	1:E:185:ARG:HD3	1.94	0.48
1:G:125:HIS:O	1:G:129:THR:HG23	2.14	0.48
1:A:62:GLU:HG3	1:A:67:LYS:CG	2.39	0.48
1:F:326:ILE:O	1:F:330:HIS:ND1	2.39	0.48
1:G:129:THR:OG1	1:G:130:HIS:N	2.32	0.48
1:B:150:TRP:CD2	1:B:265:LYS:HE3	2.49	0.48
1:D:231:PRO:HG2	1:D:248:LEU:HD13	1.96	0.48
1:B:59:LYS:O	1:B:62:GLU:HG2	2.13	0.48
1:E:283:ASN:ND2	1:E:301:TYR:CZ	2.82	0.48
1:A:69:ASN:O	1:A:73:GLU:OE2	2.32	0.47
1:A:118:LYS:HG3	1:A:318:VAL:HG13	1.96	0.47
1:E:73:GLU:OE2	1:E:74:LYS:NZ	2.47	0.47
1:E:139:TYR:CE1	1:E:273:ILE:HD12	2.49	0.47
1:F:87:PHE:HZ	1:G:95:GLU:HB3	1.79	0.47
1:F:276:TRP:CD1	1:F:327:GLU:HB3	2.49	0.47
1:D:15:PHE:CE2	1:D:336:LEU:HD22	2.50	0.47
1:D:131:GLY:O	1:D:133:ASP:N	2.47	0.47
1:F:190:LYS:HG3	1:F:191:SER:H	1.79	0.47
1:B:223:MET:HE2	1:C:99:ARG:NH1	2.29	0.47
1:C:277:TYR:OH	1:C:283:ASN:HB3	2.15	0.47
1:B:252:GLY:HA3	1:B:257:ALA:HA	1.96	0.47
1:A:10:ALA:O	1:A:11:SER:HB3	2.15	0.47
1:A:70:ALA:O	1:A:73:GLU:CD	2.58	0.47
1:C:50:VAL:HG22	1:D:264:GLN:NE2	2.30	0.47
1:C:185:ARG:HD3	1:C:185:ARG:N	2.30	0.47
1:D:34:LYS:HG2	1:D:220:PHE:CD2	2.49	0.47
1:A:134:ILE:H	1:A:134:ILE:CD1	2.14	0.47
1:C:134:ILE:HD13	1:C:330:HIS:CE1	2.49	0.47
1:G:166:LEU:HB2	1:G:214:LEU:HD23	1.96	0.47
1:B:164:THR:HG22	1:B:173:LEU:HA	1.97	0.47
1:E:69:ASN:ND2	1:F:299:VAL:HG23	2.29	0.47
1:F:28:SER:O	1:F:40:TRP:HA	2.15	0.47
1:F:292:GLY:HA3	1:F:302:ARG:HB2	1.97	0.47
1:C:12:VAL:O	1:C:111:ASN:N	2.46	0.47
1:C:139:TYR:CE1	1:C:273:ILE:CD1	2.98	0.47
1:D:165:ARG:HE	1:D:202:LEU:HD11	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:185:ARG:HG2	1:F:186:GLN:NE2	2.30	0.47
1:D:196:LEU:HD23	1:D:196:LEU:O	2.15	0.46
1:E:13:LEU:HD23	1:E:15:PHE:CE2	2.50	0.46
1:C:145:ASN:O	1:C:179:ALA:HB1	2.14	0.46
1:C:302:ARG:HD3	1:C:308:LYS:O	2.15	0.46
1:F:72:ILE:HD13	1:G:293:SER:HB3	1.96	0.46
1:G:325:THR:HG23	1:G:328:ASP:HB2	1.98	0.46
1:A:60:LYS:O	1:A:62:GLU:HA	2.15	0.46
1:A:231:PRO:HD2	1:A:248:LEU:HD13	1.98	0.46
1:A:244:LYS:HD2	1:A:247:PHE:CE2	2.51	0.46
1:B:127:ILE:O	1:B:131:GLY:HA2	2.15	0.46
1:B:262:HIS:HB3	1:B:264:GLN:OE1	2.16	0.46
1:C:26:PHE:CE1	1:C:261:ILE:HG13	2.51	0.46
1:D:57:ARG:HG2	1:E:311:TYR:CD2	2.50	0.46
1:B:309:ASP:O	1:B:313:LEU:HG	2.15	0.46
1:E:279:ASP:OD1	1:E:279:ASP:N	2.49	0.46
1:B:277:TYR:OH	1:B:283:ASN:HB3	2.14	0.46
1:B:297:GLN:OE1	1:B:301:TYR:OH	2.33	0.46
1:C:304:PRO:HA	1:C:309:ASP:OD1	2.16	0.46
1:D:243:THR:HG23	1:D:244:LYS:H	1.80	0.46
1:E:24:ALA:HB1	1:E:94:VAL:CG1	2.46	0.46
1:E:277:TYR:OH	1:E:283:ASN:HB3	2.15	0.46
1:B:24:ALA:HB1	1:B:94:VAL:HG13	1.97	0.46
1:B:286:ILE:HD13	1:B:301:TYR:HD2	1.81	0.46
1:D:131:GLY:C	1:D:133:ASP:H	2.23	0.46
1:E:115:TYR:CE1	1:E:119:LEU:HD13	2.51	0.46
1:B:274:ASP:HB2	1:B:288:VAL:HG22	1.98	0.46
1:E:59:LYS:HB2	1:E:62:GLU:HG2	1.98	0.46
1:A:280:ALA:O	1:A:281:GLU:HB2	2.16	0.46
1:A:289:GLU:OE1	1:A:338:ARG:NH2	2.43	0.46
1:E:140:ALA:CB	1:E:199:LEU:HD23	2.43	0.46
1:E:165:ARG:HD2	1:E:202:LEU:HD21	1.97	0.46
1:D:126:TYR:CD2	1:D:127:ILE:HD12	2.47	0.46
1:D:136:GLY:HA3	1:D:204:THR:CG2	2.45	0.46
1:B:119:LEU:HD13	1:B:318:VAL:HG21	1.97	0.45
1:E:72:ILE:HD11	1:F:300:ALA:HB2	1.98	0.45
1:C:73:GLU:OE1	1:E:111:ASN:OD1	2.34	0.45
1:F:274:ASP:CG	1:F:302:ARG:HH21	2.24	0.45
1:G:74:LYS:HD2	1:G:75:PRO:HD2	0.69	0.45
1:A:244:LYS:HD2	1:A:247:PHE:CD2	2.51	0.45
1:D:113:MET:HA	1:D:113:MET:HE2	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:51:ARG:HH12	1:G:262:HIS:CD2	2.31	0.45
1:B:11:SER:HB3	1:B:115:TYR:CE2	2.52	0.45
1:C:268:ASN:HA	1:C:271:ARG:CZ	2.47	0.45
1:E:59:LYS:HD2	1:E:62:GLU:CG	2.47	0.45
1:E:72:ILE:HD13	1:F:293:SER:CB	2.45	0.45
1:G:139:TYR:O	1:G:140:ALA:C	2.59	0.45
1:D:28:SER:HB2	1:D:85:LEU:HD23	1.98	0.45
1:D:72:ILE:HD12	1:E:293:SER:HB3	1.97	0.45
1:E:126:TYR:CD1	1:E:126:TYR:C	2.95	0.45
1:F:163:ILE:HD12	1:F:199:LEU:HD11	1.97	0.45
1:F:331:PHE:O	1:F:335:VAL:HG23	2.16	0.45
1:A:25:LEU:HD13	1:A:258:ILE:HB	1.99	0.45
1:C:23:ASP:OD2	1:C:249:TYR:OH	2.32	0.45
1:C:237:LEU:HD12	1:C:238:ASP:HB2	1.99	0.45
1:F:153:ARG:HG2	1:F:159:ILE:CD1	2.47	0.45
1:B:19:LEU:HD11	1:B:139:TYR:CD2	2.52	0.45
1:D:125:HIS:O	1:D:129:THR:OG1	2.26	0.45
1:F:276:TRP:HB2	1:F:302:ARG:HH22	1.82	0.45
1:A:70:ALA:CA	1:A:73:GLU:OE2	2.65	0.45
1:A:276:TRP:HB3	1:A:327:GLU:HB3	1.99	0.45
1:B:150:TRP:CE2	1:B:265:LYS:HE3	2.52	0.45
1:D:262:HIS:HB3	1:D:264:GLN:CD	2.42	0.45
1:E:138:ARG:CD	1:E:273:ILE:HG23	2.47	0.45
1:F:140:ALA:HB1	1:F:199:LEU:HD22	1.99	0.45
1:F:145:ASN:O	1:F:153:ARG:NH1	2.47	0.45
1:F:306:ALA:O	1:F:308:LYS:HG2	2.17	0.45
1:A:143:LEU:HB3	1:A:163:ILE:HD11	1.99	0.44
1:A:301:TYR:O	1:A:302:ARG:HB2	2.18	0.44
1:B:74:LYS:HB3	1:B:74:LYS:HE3	1.78	0.44
1:D:149:LEU:HB2	1:D:153:ARG:HG3	1.98	0.44
1:E:80:VAL:HG12	1:E:245:SER:HB3	1.99	0.44
1:F:100:VAL:HB	1:F:213:VAL:HG21	1.98	0.44
1:F:185:ARG:HG2	1:F:186:GLN:HE21	1.82	0.44
1:F:231:PRO:CG	1:F:248:LEU:HD13	2.47	0.44
1:D:138:ARG:HB3	1:D:273:ILE:CG1	2.47	0.44
1:A:134:ILE:HD13	1:A:330:HIS:CE1	2.52	0.44
1:B:282:ILE:O	1:B:284:GLY:N	2.46	0.44
1:D:62:GLU:HA	1:D:67:LYS:HE3	1.98	0.44
1:D:166:LEU:HB2	1:D:214:LEU:HD23	1.98	0.44
1:E:249:TYR:CD2	1:E:260:ALA:HB3	2.52	0.44
1:E:301:TYR:H	1:E:301:TYR:HD1	1.63	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:68:LEU:HD12	1:G:304:PRO:CD	2.47	0.44
1:F:249:TYR:CD2	1:F:260:ALA:HB3	2.52	0.44
1:B:153:ARG:HG2	1:B:159:ILE:CD1	2.46	0.44
1:D:50:VAL:HG22	1:E:264:GLN:NE2	2.32	0.44
1:D:232:SER:CB	1:D:262:HIS:CD2	3.01	0.44
1:G:242:SER:O	1:G:243:THR:HG22	2.17	0.44
1:B:100:VAL:HB	1:B:213:VAL:HG21	1.99	0.44
1:D:46:ARG:HH12	1:D:84:THR:CG2	2.30	0.44
1:D:252:GLY:O	1:D:257:ALA:HB2	2.17	0.44
1:D:277:TYR:CE1	1:D:286:ILE:HG12	2.53	0.44
1:A:139:TYR:CD1	1:A:273:ILE:CD1	3.00	0.44
1:A:146:GLY:HA2	1:A:148:PHE:CE2	2.53	0.44
1:B:137:ASN:OD1	1:B:200:ALA:HB1	2.18	0.44
1:D:249:TYR:CD2	1:D:260:ALA:HB3	2.52	0.44
1:E:201:LYS:HA	1:E:204:THR:OG1	2.17	0.44
1:G:183:PRO:HG2	1:G:186:GLN:O	2.18	0.44
1:G:297:GLN:O	1:G:299:VAL:HG23	2.17	0.44
1:C:70:ALA:O	1:C:73:GLU:HG3	2.18	0.44
1:D:137:ASN:HA	1:D:200:ALA:HB1	2.00	0.44
1:E:16:GLU:O	1:E:18:LYS:HD2	2.18	0.44
1:E:134:ILE:O	1:E:138:ARG:HG3	2.18	0.44
1:G:185:ARG:HD3	1:G:185:ARG:N	2.31	0.44
1:A:271:ARG:HG3	1:A:288:VAL:HB	2.00	0.44
1:B:26:PHE:CE1	1:B:261:ILE:HG13	2.52	0.44
1:F:26:PHE:HB3	1:F:43:ILE:HD12	1.99	0.44
1:G:75:PRO:HB3	1:G:77:LEU:HD23	1.99	0.44
1:C:150:TRP:CD2	1:C:265:LYS:HE3	2.53	0.43
1:D:173:LEU:C	1:D:173:LEU:HD13	2.43	0.43
1:F:31:TRP:CE3	1:F:222:ARG:HB2	2.53	0.43
1:F:116:ARG:O	1:F:120:VAL:HG23	2.18	0.43
1:F:125:HIS:O	1:F:129:THR:HG23	2.18	0.43
1:F:325:THR:HG23	1:F:328:ASP:HB2	2.00	0.43
1:A:12:VAL:O	1:A:111:ASN:HB2	2.18	0.43
1:B:62:GLU:HA	1:B:67:LYS:HE2	1.99	0.43
1:D:138:ARG:HG2	1:D:187:PHE:CD1	2.53	0.43
1:E:13:LEU:HD21	1:E:119:LEU:HD22	1.99	0.43
1:E:146:GLY:HA2	1:E:148:PHE:CE2	2.53	0.43
1:E:166:LEU:HD11	1:E:216:ARG:HB2	1.99	0.43
1:F:225:GLU:OE1	1:G:216:ARG:NH2	2.51	0.43
1:G:132:LEU:O	1:G:132:LEU:CD1	2.66	0.43
1:D:150:TRP:NE1	1:D:265:LYS:HG3	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:325:THR:HG23	1:D:328:ASP:HB2	2.00	0.43
1:E:122:THR:HG22	1:E:317:TRP:CE2	2.53	0.43
1:E:250:SER:HA	1:E:258:ILE:O	2.18	0.43
1:B:279:ASP:CG	1:B:282:ILE:HD11	2.43	0.43
1:E:118:LYS:HB3	1:E:318:VAL:HG13	2.00	0.43
1:B:115:TYR:HE1	1:B:119:LEU:HD22	1.83	0.43
1:B:277:TYR:HB2	1:B:278:PRO:HD2	1.99	0.43
1:E:134:ILE:HA	1:E:137:ASN:HB2	1.99	0.43
1:E:309:ASP:O	1:E:313:LEU:HG	2.18	0.43
1:B:12:VAL:HB	1:B:111:ASN:CB	2.40	0.43
1:B:100:VAL:HB	1:B:213:VAL:CG2	2.49	0.43
1:C:149:LEU:HD13	1:C:221:ILE:HD11	2.00	0.43
1:F:192:VAL:O	1:F:193:SER:HB3	2.18	0.43
1:D:144:ALA:O	1:D:179:ALA:HA	2.18	0.43
1:B:153:ARG:NH1	1:B:184:LEU:HD11	2.33	0.43
1:D:27:PHE:O	1:D:92:LEU:HD12	2.19	0.43
1:D:46:ARG:HH12	1:D:84:THR:HG23	1.83	0.43
1:F:304:PRO:HA	1:F:309:ASP:OD2	2.18	0.43
1:B:249:TYR:CG	1:B:262:HIS:CE1	3.06	0.43
1:D:57:ARG:HG2	1:E:311:TYR:CE2	2.53	0.43
1:E:18:LYS:HD2	1:E:18:LYS:N	2.33	0.43
1:F:65:PRO:CA	1:G:303:GLN:HE22	2.27	0.43
1:G:74:LYS:HD2	1:G:74:LYS:C	2.42	0.43
1:B:62:GLU:HG3	1:B:62:GLU:O	2.19	0.42
1:B:165:ARG:HA	1:B:215:LEU:HD23	2.01	0.42
1:E:163:ILE:HG21	1:E:199:LEU:CD1	2.49	0.42
1:E:315:ASP:O	1:E:319:LEU:N	2.50	0.42
1:F:78:GLN:HB3	1:F:245:SER:HB2	2.01	0.42
1:G:149:LEU:HD22	1:G:221:ILE:HD11	2.01	0.42
1:B:81:ASP:O	1:B:246:ARG:NH2	2.52	0.42
1:B:264:GLN:OE1	1:B:264:GLN:N	2.49	0.42
1:D:73:GLU:HG2	1:E:298:GLY:HA3	1.99	0.42
1:D:277:TYR:OH	1:D:283:ASN:ND2	2.52	0.42
1:E:118:LYS:O	1:E:122:THR:OG1	2.27	0.42
1:E:132:LEU:O	1:E:204:THR:HG22	2.18	0.42
1:D:153:ARG:CZ	1:D:184:LEU:HD11	2.49	0.42
1:B:15:PHE:CD2	1:B:336:LEU:HD22	2.54	0.42
1:D:262:HIS:O	1:D:263:SER:C	2.62	0.42
1:B:223:MET:HE2	1:C:99:ARG:HH12	1.84	0.42
1:D:326:ILE:HA	1:D:329:GLN:HB2	2.02	0.42
1:E:62:GLU:O	1:E:63:ALA:C	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:MET:N	1:B:113:MET:SD	2.91	0.42
1:D:68:LEU:HD11	1:E:291:TYR:CD2	2.54	0.42
1:E:25:LEU:HD22	1:E:260:ALA:HB2	2.00	0.42
1:E:70:ALA:O	1:E:73:GLU:HG3	2.19	0.42
1:E:144:ALA:O	1:E:179:ALA:HA	2.20	0.42
1:C:134:ILE:H	1:C:134:ILE:CD1	2.28	0.42
1:F:72:ILE:HD11	1:G:300:ALA:HB2	2.01	0.42
1:F:221:ILE:HG22	1:F:223:MET:HG3	2.01	0.42
1:B:48:LYS:HE2	1:B:82:VAL:HG11	2.02	0.42
1:C:277:TYR:CE1	1:C:286:ILE:HG12	2.55	0.42
1:E:308:LYS:HD2	1:E:328:ASP:OD2	2.20	0.42
1:G:46:ARG:HH21	1:G:84:THR:CG2	2.32	0.42
1:E:277:TYR:CD2	1:E:301:TYR:HB3	2.55	0.42
1:F:313:LEU:HD13	1:F:328:ASP:HB3	2.02	0.42
1:A:72:ILE:HD12	1:B:300:ALA:HB2	2.02	0.42
1:B:189:GLU:H	1:B:189:GLU:HG2	1.56	0.42
1:F:51:ARG:HA	1:F:77:LEU:HD22	2.01	0.42
1:A:165:ARG:HH21	1:A:202:LEU:HD11	1.85	0.41
1:A:279:ASP:O	1:A:282:ILE:HG13	2.19	0.41
1:B:138:ARG:HB3	1:B:273:ILE:CG2	2.50	0.41
1:B:277:TYR:OH	1:B:284:GLY:O	2.29	0.41
1:C:77:LEU:HD11	1:D:235:LEU:HD13	2.02	0.41
1:D:253:GLN:NE2	1:D:256:LYS:HD2	2.35	0.41
1:D:277:TYR:OH	1:D:284:GLY:O	2.29	0.41
1:D:291:TYR:HD1	1:D:309:ASP:HB2	1.85	0.41
1:E:69:ASN:HD21	1:F:299:VAL:CG2	2.32	0.41
1:E:241:LYS:O	1:E:241:LYS:HG2	2.20	0.41
1:G:132:LEU:O	1:G:204:THR:HG22	2.20	0.41
1:G:324:PRO:O	1:G:329:GLN:NE2	2.49	0.41
1:A:73:GLU:OE2	1:A:73:GLU:N	2.50	0.41
1:F:185:ARG:HD3	1:F:185:ARG:N	2.35	0.41
1:D:68:LEU:C	1:D:68:LEU:HD23	2.45	0.41
1:E:273:ILE:HD13	1:E:273:ILE:HG21	1.80	0.41
1:F:297:GLN:HA	1:G:110:CYS:O	2.21	0.41
1:F:327:GLU:H	1:F:327:GLU:CD	2.28	0.41
1:B:126:TYR:CD1	1:B:126:TYR:C	2.98	0.41
1:D:140:ALA:CB	1:D:199:LEU:HD13	2.48	0.41
1:A:61:GLY:HA2	1:A:62:GLU:CB	2.48	0.41
1:C:77:LEU:HD11	1:D:235:LEU:CD1	2.51	0.41
1:C:159:ILE:HA	1:C:220:PHE:O	2.20	0.41
1:E:149:LEU:HD23	1:E:149:LEU:HA	1.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:307:LYS:H	1:A:307:LYS:HG2	1.53	0.41
1:E:193:SER:O	1:E:197:GLU:HB2	2.21	0.41
1:A:188:GLU:H	1:A:188:GLU:HG2	1.61	0.41
1:A:177:PHE:CZ	1:A:199:LEU:HD12	2.56	0.41
1:B:135:LEU:HD23	1:B:135:LEU:HA	1.89	0.41
1:D:72:ILE:HD12	1:E:293:SER:CB	2.50	0.41
1:D:305:LYS:HE2	1:D:305:LYS:HB3	1.84	0.41
1:D:325:THR:HG23	1:D:328:ASP:H	1.85	0.41
1:E:76:ASN:C	1:E:76:ASN:HD22	2.29	0.41
1:G:134:ILE:HD12	1:G:137:ASN:HD22	1.85	0.41
1:B:151:ARG:HG2	1:B:230:PHE:CD2	2.56	0.41
1:B:166:LEU:HB2	1:B:214:LEU:HD23	2.03	0.41
1:G:112:SER:OG	1:G:115:TYR:HB2	2.21	0.41
1:A:34:LYS:HG2	1:A:220:PHE:CG	2.56	0.40
1:A:139:TYR:CD1	1:A:273:ILE:HD12	2.56	0.40
1:B:148:PHE:CD1	1:B:148:PHE:C	3.00	0.40
1:B:149:LEU:HD22	1:B:221:ILE:HD11	2.03	0.40
1:D:118:LYS:O	1:D:122:THR:HG23	2.21	0.40
1:A:70:ALA:HA	1:A:73:GLU:CD	2.46	0.40
1:B:41:GLN:CD	1:B:42:PRO:HD2	2.45	0.40
1:B:173:LEU:HD13	1:B:175:GLY:N	2.36	0.40
1:B:221:ILE:HD12	1:B:221:ILE:N	2.36	0.40
1:C:27:PHE:HB2	1:C:93:LYS:HB3	2.04	0.40
1:D:31:TRP:CE3	1:D:222:ARG:HB2	2.56	0.40
1:D:222:ARG:NH1	1:E:168:GLY:HA2	2.36	0.40
1:E:159:ILE:HA	1:E:220:PHE:O	2.21	0.40
1:E:162:GLN:HG2	1:E:176:VAL:HG22	2.03	0.40
1:A:308:LYS:HE3	1:A:308:LYS:HB3	1.97	0.40
1:B:17:ARG:NE	1:B:20:ASP:OD1	2.54	0.40
1:B:30:ASN:HA	1:B:90:ASP:OD1	2.22	0.40
1:D:191:SER:O	1:D:191:SER:OG	2.35	0.40
1:E:307:LYS:H	1:E:307:LYS:HG2	1.64	0.40
1:G:86:PRO:HB2	1:G:89:SER:HB2	2.03	0.40
1:G:189:GLU:O	1:G:191:SER:N	2.54	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:138:ARG:NH2	1:D:281:GLU:OE1[1_565]	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	331/348 (95%)	307 (93%)	15 (4%)	9 (3%)	4	16
1	B	331/348 (95%)	309 (93%)	18 (5%)	4 (1%)	10	34
1	C	331/348 (95%)	301 (91%)	26 (8%)	4 (1%)	10	34
1	D	331/348 (95%)	302 (91%)	21 (6%)	8 (2%)	4	18
1	E	332/348 (95%)	308 (93%)	17 (5%)	7 (2%)	5	21
1	F	332/348 (95%)	312 (94%)	17 (5%)	3 (1%)	14	41
1	G	332/348 (95%)	304 (92%)	22 (7%)	6 (2%)	6	25
All	All	2320/2436 (95%)	2143 (92%)	136 (6%)	41 (2%)	6	25

All (41) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	190	LYS
1	A	281	GLU
1	A	307	LYS
1	B	192	VAL
1	B	280	ALA
1	B	307	LYS
1	C	307	LYS
1	D	129	THR
1	D	193	SER
1	D	307	LYS
1	E	63	ALA
1	E	307	LYS
1	F	129	THR
1	F	193	SER
1	G	60	LYS
1	A	302	ARG
1	C	150	TRP
1	C	281	GLU

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Mol	Chain	Res	Type
1	D	62	GLU
1	D	246	ARG
1	E	129	THR
1	E	132	LEU
1	E	192	VAL
1	E	274	ASP
1	F	190	LYS
1	G	75	PRO
1	G	129	THR
1	A	75	PRO
1	D	278	PRO
1	G	253	GLN
1	A	150	TRP
1	D	132	LEU
1	G	190	LYS
1	A	132	LEU
1	A	273	ILE
1	C	273	ILE
1	D	150	TRP
1	E	253	GLN
1	B	132	LEU
1	G	305	LYS
1	A	252	GLY

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	274/286 (96%)	274 (100%)	0	100	100
1	B	274/286 (96%)	273 (100%)	1 (0%)	84	94
1	C	274/286 (96%)	271 (99%)	3 (1%)	65	88
1	D	274/286 (96%)	271 (99%)	3 (1%)	65	88
1	E	274/286 (96%)	273 (100%)	1 (0%)	84	94
1	F	274/286 (96%)	269 (98%)	5 (2%)	51	80

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	274/286 (96%)	273 (100%)	1 (0%)	84	94
All	All	1918/2002 (96%)	1904 (99%)	14 (1%)	76	92

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	223	MET
1	C	49	SER
1	C	185	ARG
1	C	254	ASP
1	D	199	LEU
1	D	233	GLN
1	D	296	THR
1	E	104	VAL
1	F	82	VAL
1	F	132	LEU
1	F	165	ARG
1	F	189	GLU
1	F	281	GLU
1	G	186	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	ASN
1	A	76	ASN
1	A	268	ASN
1	A	283	ASN
1	B	262	HIS
1	C	227	GLN
1	D	262	HIS
1	D	268	ASN
1	E	76	ASN
1	E	111	ASN
1	E	130	HIS
1	E	145	ASN
1	F	33	ASN
1	F	125	HIS
1	F	186	GLN
1	F	227	GLN
1	F	233	GLN

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Mol	Chain	Res	Type
1	G	130	HIS
1	G	210	GLN
1	G	253	GLN
1	G	262	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	333/348 (95%)	0.02	10 (3%) 52 43	46, 66, 122, 160	0
1	B	333/348 (95%)	0.21	9 (2%) 56 47	45, 78, 126, 167	0
1	C	333/348 (95%)	-0.08	7 (2%) 63 54	43, 66, 108, 156	0
1	D	333/348 (95%)	0.06	15 (4%) 38 30	43, 63, 125, 179	0
1	E	334/348 (95%)	0.24	18 (5%) 31 24	45, 77, 127, 171	0
1	F	334/348 (95%)	0.04	10 (2%) 52 43	47, 70, 131, 164	0
1	G	334/348 (95%)	0.22	14 (4%) 40 32	48, 79, 128, 199	0
All	All	2334/2436 (95%)	0.10	83 (3%) 46 38	43, 72, 126, 199	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	12	VAL	5.6
1	E	12	VAL	5.5
1	C	237	LEU	5.3
1	G	12	VAL	5.2
1	F	10	ALA	5.1
1	B	12	VAL	4.8
1	G	192	VAL	4.6
1	G	10	ALA	4.6
1	G	243	THR	4.5
1	C	253	GLN	4.4
1	A	243	THR	4.2
1	D	243	THR	4.1
1	A	12	VAL	3.9
1	F	12	VAL	3.9
1	F	243	THR	3.9
1	E	237	LEU	3.9
1	E	343	GLY	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	192	VAL	3.5
1	A	237	LEU	3.5
1	E	63	ALA	3.4
1	D	241	LYS	3.4
1	D	10	ALA	3.3
1	B	243	THR	3.3
1	D	245	SER	3.3
1	E	319	LEU	3.3
1	C	10	ALA	3.2
1	B	237	LEU	3.1
1	C	243	THR	3.1
1	E	192	VAL	3.0
1	G	11	SER	2.9
1	E	239	LYS	2.9
1	F	238	ASP	2.9
1	A	10	ALA	2.9
1	B	239	LYS	2.9
1	G	130	HIS	2.9
1	D	237	LEU	2.9
1	G	319	LEU	2.9
1	C	12	VAL	2.8
1	F	192	VAL	2.7
1	E	111	ASN	2.7
1	D	236	LEU	2.7
1	G	239	LYS	2.7
1	F	253	GLN	2.7
1	F	11	SER	2.7
1	E	10	ALA	2.6
1	E	130	HIS	2.6
1	A	254	ASP	2.6
1	A	242	SER	2.6
1	D	191	SER	2.6
1	B	253	GLN	2.6
1	D	51	ARG	2.5
1	A	241	LYS	2.5
1	F	242	SER	2.5
1	E	241	LYS	2.5
1	D	63	ALA	2.5
1	E	134	ILE	2.4
1	A	63	ALA	2.4
1	G	241	LYS	2.4
1	D	192	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	113	MET	2.4
1	E	236	LEU	2.4
1	D	239	LYS	2.4
1	G	343	GLY	2.4
1	F	241	LYS	2.3
1	F	63	ALA	2.3
1	G	237	LEU	2.3
1	C	241	LYS	2.2
1	D	242	SER	2.2
1	G	308	LYS	2.2
1	G	191	SER	2.2
1	E	299	VAL	2.1
1	B	241	LYS	2.1
1	C	239	LYS	2.1
1	A	11	SER	2.1
1	E	295	THR	2.1
1	D	189	GLU	2.1
1	E	253	GLN	2.1
1	A	253	GLN	2.0
1	B	130	HIS	2.0
1	E	306	ALA	2.0
1	B	191	SER	2.0
1	E	288	VAL	2.0
1	G	255	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.