



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:55 PM GMT

PDB ID : 4JTV
Title : Crystal structure of 2009 pandemic influenza virus hemagglutinin complexed with human receptor analogue LSTc
Authors : Zhang, W.; Shi, Y.; Qi, J.; Gao, F.; Li, Q.; Fan, Z.; Yan, J.; Gao, G.F.
Deposited on : 2013-03-24
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

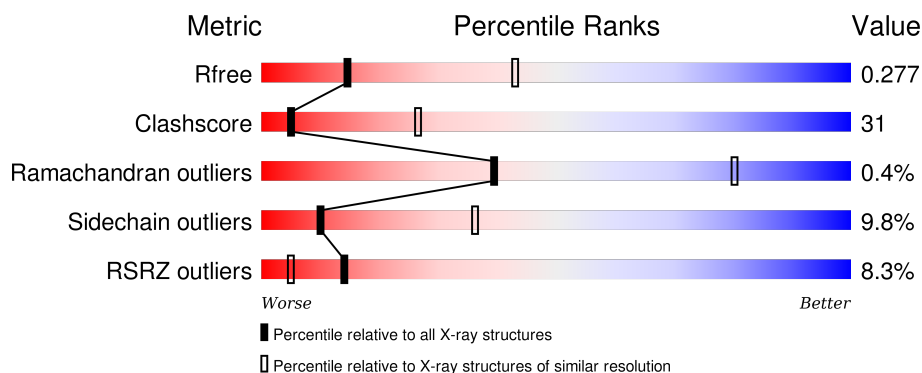
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	321	<div> <div>4%</div> <div>48% 45% 7%</div> </div>
1	C	321	<div> <div>7%</div> <div>52% 40% 7%</div> </div>
1	E	321	<div> <div>3%</div> <div>56% 38% 5%</div> </div>
1	G	321	<div> <div>6%</div> <div>60% 36%</div> </div>
1	I	321	<div> <div>4%</div> <div>52% 40% 7%</div> </div>

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Mol	Chain	Length	Quality of chain
1	K	321	
2	B	162	
2	D	162	
2	F	162	
2	H	162	
2	J	162	
2	L	162	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAG	G	602	-	-	-	X

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 23528 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 Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	321	Total	C	N	O	S	0	0	0
			2505	1584	433	477	11			
1	C	321	Total	C	N	O	S	0	0	0
			2509	1586	433	479	11			
1	E	321	Total	C	N	O	S	0	0	0
			2509	1586	433	479	11			
1	G	321	Total	C	N	O	S	0	0	0
			2505	1584	433	477	11			
1	I	321	Total	C	N	O	S	0	0	0
			2510	1587	433	479	11			
1	K	321	Total	C	N	O	S	0	0	0
			2509	1586	433	479	11			

- Molecule 2 is a protein called Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	162	Total	C	N	O	S	0	0	0
			1305	822	220	257	6			
2	D	162	Total	C	N	O	S	0	0	0
			1300	818	219	257	6			
2	F	161	Total	C	N	O	S	0	0	0
			1302	821	219	256	6			
2	H	162	Total	C	N	O	S	0	0	0
			1305	822	220	257	6			
2	J	162	Total	C	N	O	S	0	0	0
			1300	818	219	257	6			
2	L	161	Total	C	N	O	S	0	0	0
			1302	821	219	256	6			

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	G	1	Total	C	N	O	0	0
			14	8	1	5		
3	G	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	3	Total	C	N	O	0	0
			45	25	2	18		
4	I	3	Total	C	N	O	0	0
			45	25	2	18		
4	K	3	Total	C	N	O	0	0
			45	25	2	18		

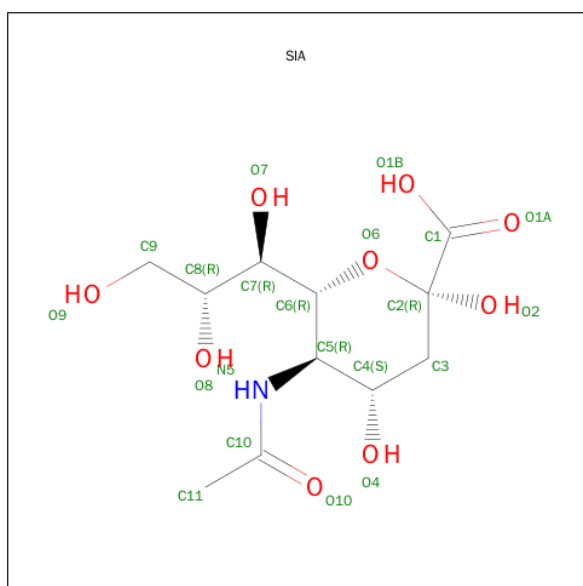
- Molecule 5 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	C	4	Total	C	N	O	0	0
			56	31	2	23		
5	E	4	Total	C	N	O	0	0
			56	31	2	23		

- Molecule 6 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	E	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 7 is SUGAR (O-SIALIC ACID) (three-letter code: SIA) (formula: $C_{11}H_{19}NO_9$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	G	1	Total	C	N	O	0	0
			20	11	1	8		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	40	Total	O	0	0
			40	40		
8	B	13	Total	O	0	0
			13	13		
8	C	29	Total	O	0	0
			29	29		

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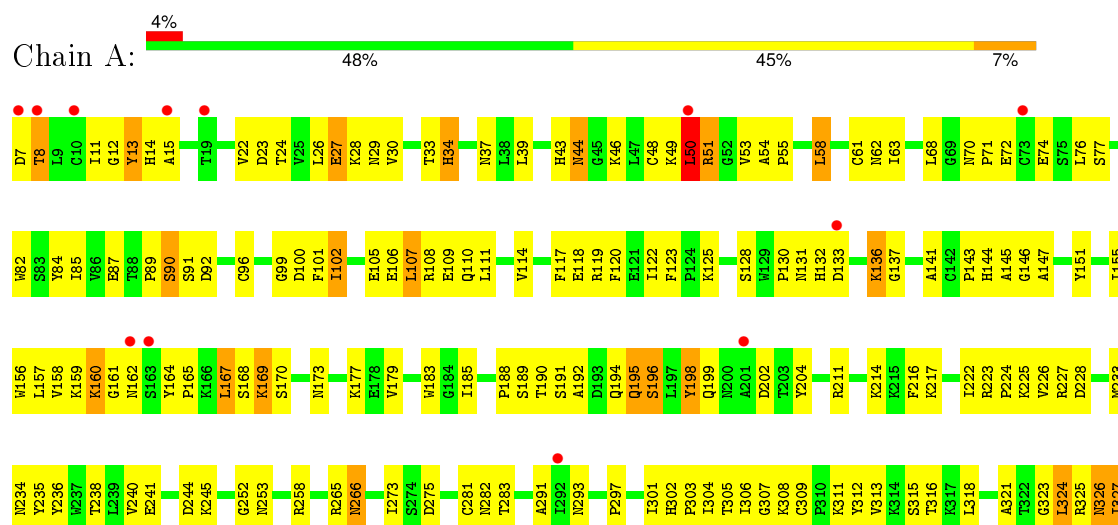
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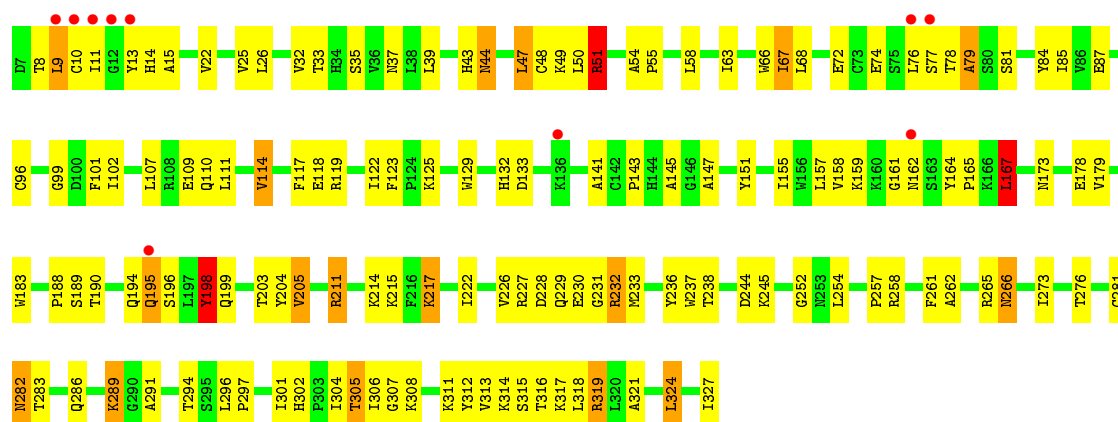
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	D	13	Total 13	O 13	0	0
8	E	31	Total 31	O 31	0	0
8	F	7	Total 7	O 7	0	0
8	G	19	Total 19	O 19	0	0
8	H	13	Total 13	O 13	0	0
8	I	35	Total 35	O 35	0	0
8	J	18	Total 18	O 18	0	0
8	K	25	Total 25	O 25	0	0
8	L	31	Total 31	O 31	0	0

3 Residue-property plots

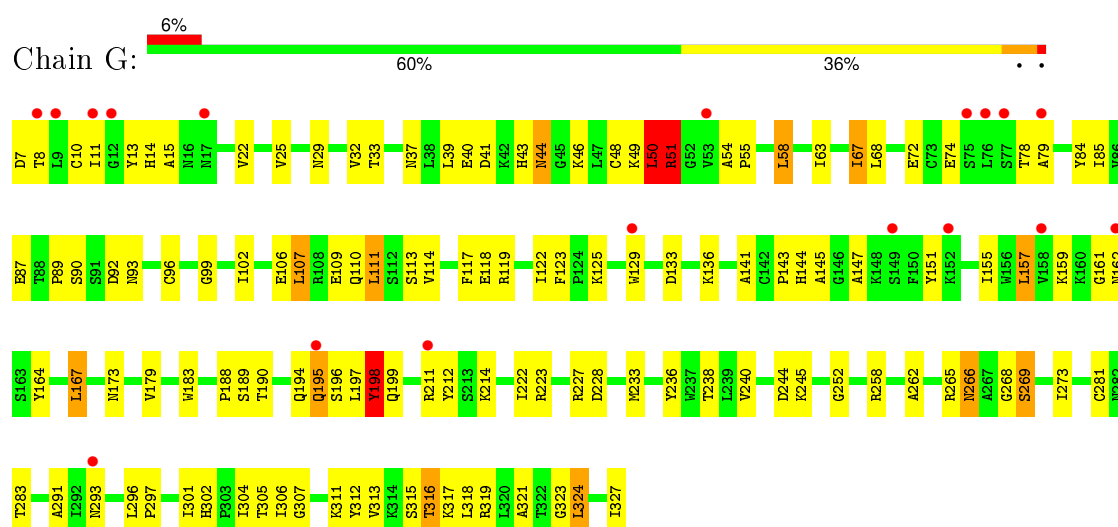
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Hemagglutinin

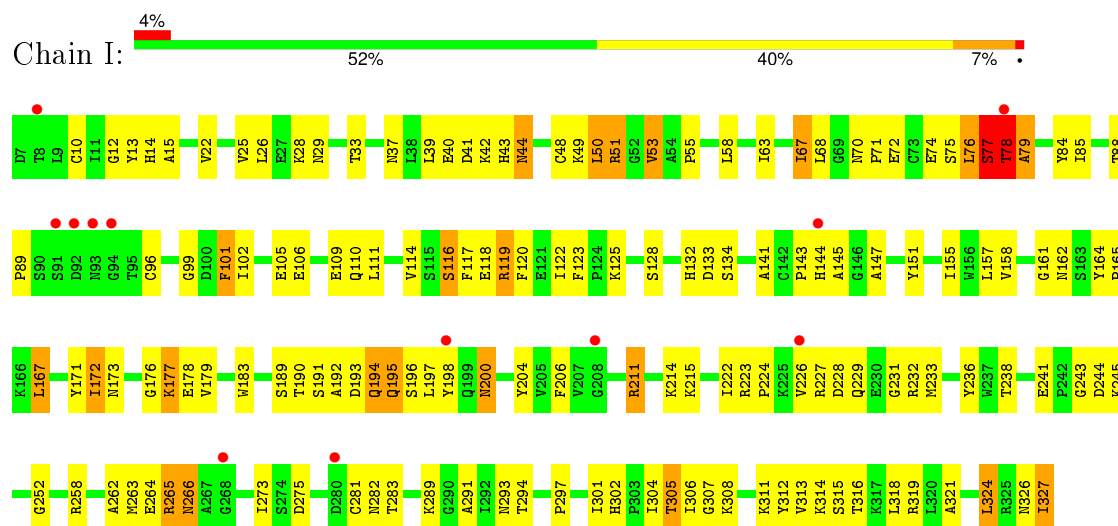




• Molecule 1: Hemagglutinin

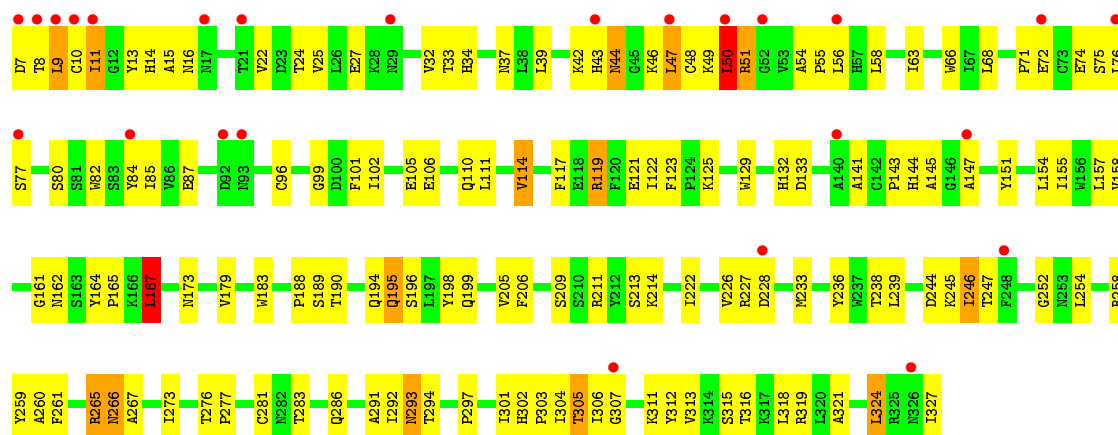


• Molecule 1: Hemagglutinin

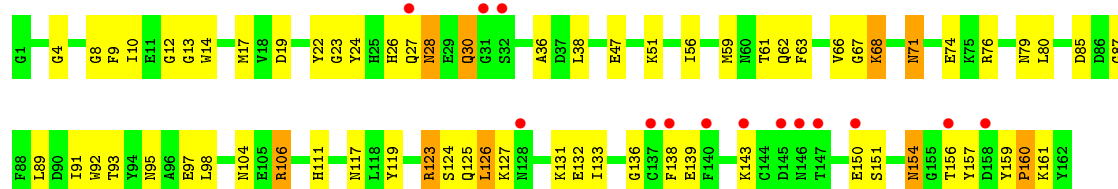


• Molecule 1: Hemagglutinin

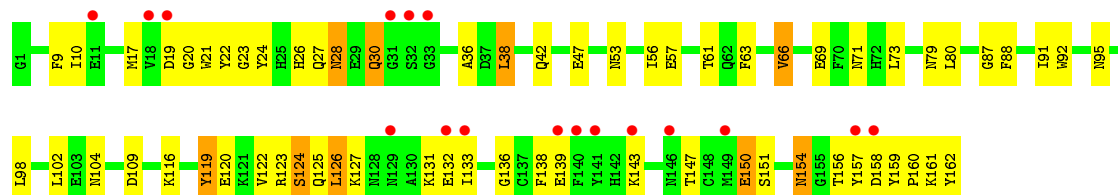




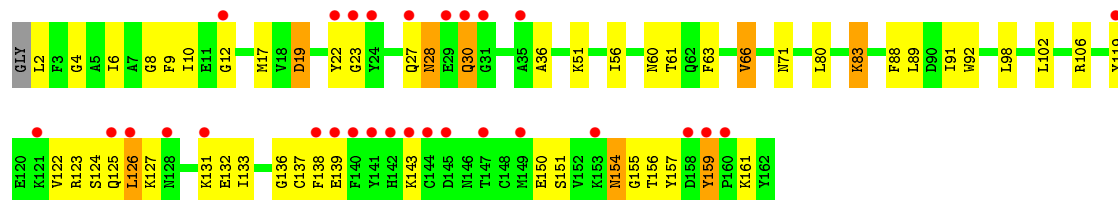
• Molecule 2: Hemagglutinin



• Molecule 2: Hemagglutinin

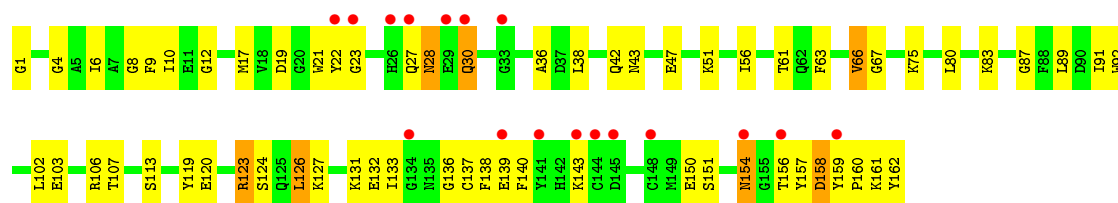


• Molecule 2: Hemagglutinin

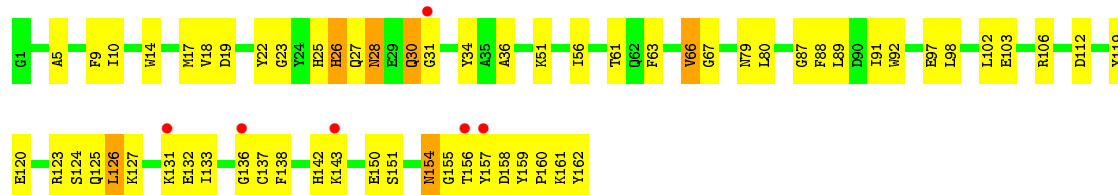


• Molecule 2: Hemagglutinin

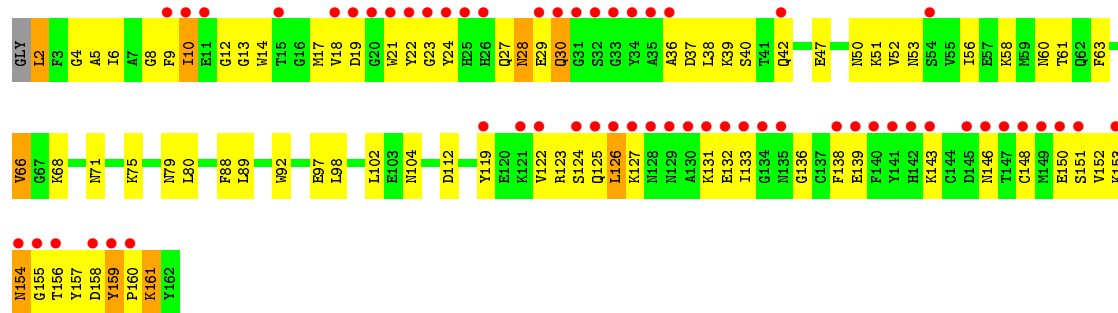




• Molecule 2: Hemagglutinin



• Molecule 2: Hemagglutinin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	66.72Å 117.32Å 117.39Å 61.78° 81.82° 77.42°	Depositor
Resolution (Å)	38.39 – 3.00 49.68 – 3.00	Depositor EDS
% Data completeness (in resolution range)	81.5 (38.39-3.00) 80.6 (49.68-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.224 , 0.275 0.227 , 0.277	Depositor DCC
R_{free} test set	2707 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	55.4	Xtriage
Anisotropy	0.701	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 58.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 53766 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	23528	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.66% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SIA, GAL, NAG

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	0.57	0/2568	0.75	5/3488 (0.1%)
1	C	0.55	0/2572	0.69	8/3493 (0.2%)
1	E	0.39	0/2572	0.58	4/3493 (0.1%)
1	G	0.33	0/2568	0.77	9/3488 (0.3%)
1	I	0.47	0/2573	0.63	3/3495 (0.1%)
1	K	0.33	0/2572	0.63	5/3493 (0.1%)
2	B	0.48	0/1333	0.54	1/1797 (0.1%)
2	D	0.38	0/1328	0.45	0/1791
2	F	0.26	0/1330	0.46	0/1794
2	H	0.26	0/1333	0.49	1/1797 (0.1%)
2	J	0.43	0/1328	0.48	0/1791
2	L	0.26	0/1330	0.53	1/1794 (0.1%)
All	All	0.42	0/23407	0.62	37/31714 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	I	0	1

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	119	ARG	NE-CZ-NH1	-16.56	112.02	120.30
1	A	119	ARG	NE-CZ-NH2	16.50	128.55	120.30
1	A	119	ARG	NE-CZ-NH1	-16.32	112.14	120.30
1	G	119	ARG	NE-CZ-NH2	15.97	128.29	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	212	TYR	N-CA-CB	11.91	132.04	110.60
1	G	211	ARG	CB-CA-C	-11.01	88.39	110.40
1	I	78	THR	N-CA-CB	-10.87	89.66	110.30
2	L	10	ILE	CB-CA-C	10.73	133.07	111.60
1	C	74	GLU	CB-CA-C	-10.72	88.96	110.40
1	C	78	THR	N-CA-CB	-9.71	91.85	110.30
1	G	50	LEU	CB-CA-C	-8.91	93.28	110.20
1	A	50	LEU	CB-CA-C	-8.80	93.47	110.20
1	C	50	LEU	CB-CA-C	-8.59	93.88	110.20
1	C	75	SER	N-CA-C	-8.45	88.19	111.00
1	I	119	ARG	NE-CZ-NH2	-8.39	116.10	120.30
1	K	119	ARG	NE-CZ-NH2	-8.26	116.17	120.30
1	A	119	ARG	CD-NE-CZ	8.15	135.01	123.60
1	G	119	ARG	CD-NE-CZ	8.06	134.89	123.60
1	K	119	ARG	NE-CZ-NH1	7.52	124.06	120.30
1	I	119	ARG	NE-CZ-NH1	7.34	123.97	120.30
1	C	198	TYR	N-CA-C	-7.07	91.92	111.00
1	E	51	ARG	N-CA-C	-7.04	91.98	111.00
1	C	51	ARG	N-CA-C	-6.90	92.38	111.00
1	C	212	TYR	CB-CA-C	-6.32	97.76	110.40
1	A	51	ARG	N-CA-C	-6.16	94.36	111.00
1	K	167	LEU	CA-CB-CG	5.94	128.97	115.30
1	E	167	LEU	CA-CB-CG	5.83	128.70	115.30
1	G	212	TYR	N-CA-C	-5.81	95.31	111.00
1	G	51	ARG	N-CA-C	-5.80	95.33	111.00
2	B	66	VAL	CB-CA-C	-5.77	100.43	111.40
1	G	198	TYR	N-CA-C	-5.77	95.42	111.00
2	H	66	VAL	CB-CA-C	-5.69	100.59	111.40
1	E	167	LEU	CB-CG-CD2	5.34	120.07	111.00
1	C	50	LEU	N-CA-C	5.18	125.00	111.00
1	K	167	LEU	CB-CG-CD2	5.10	119.67	111.00
1	E	198	TYR	N-CA-C	-5.04	97.38	111.00
1	K	50	LEU	CB-CA-C	-5.00	100.70	110.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	I	77	SER	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	2505	0	2444	227	0
1	C	2509	0	2451	186	0
1	E	2509	0	2451	159	4
1	G	2505	0	2445	137	0
1	I	2510	0	2456	183	0
1	K	2509	0	2452	188	0
2	B	1305	0	1228	97	0
2	D	1300	0	1216	81	0
2	F	1302	0	1226	65	0
2	H	1305	0	1228	73	0
2	J	1300	0	1216	89	0
2	L	1302	0	1226	126	4
3	A	56	0	52	6	0
3	C	14	0	13	2	0
3	G	28	0	26	5	0
4	A	45	0	38	4	0
4	I	45	0	38	3	0
4	K	45	0	38	0	0
5	C	56	0	47	1	0
5	E	56	0	47	2	0
6	E	28	0	25	3	0
7	G	20	0	17	1	0
8	A	40	0	0	59	0
8	B	13	0	0	11	0
8	C	29	0	0	34	0
8	D	13	0	0	6	0
8	E	31	0	0	27	0
8	F	7	0	0	6	0
8	G	19	0	0	22	0
8	H	13	0	0	22	0
8	I	35	0	0	32	0
8	J	18	0	0	22	0
8	K	25	0	0	47	0
8	L	31	0	0	48	0
All	All	23528	0	22380	1414	4

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 31.

All (1414) 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:C:11:ILE:HD13	2:D:24:TYR:CD2	1.53	1.42
1:I:190:THR:HG21	1:I:193:ASP:OD1	1.24	1.35
1:K:106:GLU:HB3	8:K:915:HOH:O	1.20	1.34
2:L:19:ASP:HA	8:L:228:HOH:O	1.33	1.28
1:A:8:THR:HG22	2:B:138:PHE:O	1.18	1.27
1:A:8:THR:CG2	2:B:138:PHE:O	1.86	1.24
1:E:281:CYS:SG	8:E:709:HOH:O	1.92	1.23
2:L:56:ILE:HB	8:L:214:HOH:O	1.32	1.23
1:G:7:ASP:N	2:H:140:PHE:H	1.36	1.22
1:A:7:ASP:O	2:B:27:GLN:O	1.58	1.22
2:B:97:GLU:CB	8:B:213:HOH:O	1.88	1.22
1:K:122:ILE:HG22	8:K:914:HOH:O	1.40	1.20
1:C:258:ARG:HB3	8:C:717:HOH:O	1.34	1.20
1:I:145:ALA:HB3	8:I:903:HOH:O	1.44	1.17
1:I:190:THR:CG2	1:I:193:ASP:OD1	1.92	1.17
1:A:326:ASN:O	1:A:327:ILE:HG12	1.43	1.17
1:C:11:ILE:CD1	2:D:24:TYR:CD2	2.28	1.16
1:I:190:THR:HG23	1:I:193:ASP:N	1.60	1.16
1:A:305:THR:CG2	1:A:309:CYS:SG	2.33	1.16
1:E:178:GLU:OE1	1:E:265:ARG:NH1	1.77	1.16
1:C:11:ILE:CD1	2:D:24:TYR:CE2	2.27	1.15
1:A:108:ARG:HB2	8:A:734:HOH:O	0.98	1.15
1:K:14:HIS:CG	1:K:15:ALA:H	1.65	1.15
1:A:326:ASN:C	1:A:327:ILE:HG12	1.59	1.14
1:E:203:THR:O	1:E:217:LYS:HE2	1.46	1.13
2:L:38:LEU:C	8:L:207:HOH:O	1.86	1.12
1:E:66:TRP:HE1	1:E:77:SER:HB2	1.02	1.12
2:H:67:GLY:N	8:H:205:HOH:O	1.82	1.12
2:B:97:GLU:HB2	8:B:213:HOH:O	1.46	1.11
1:I:88:THR:HB	8:I:908:HOH:O	1.49	1.11
1:I:190:THR:CG2	1:I:193:ASP:H	1.62	1.11
1:K:14:HIS:CE1	1:K:15:ALA:O	2.04	1.11
1:K:162:ASN:OD1	1:K:199:GLN:NE2	1.81	1.10
2:L:39:LYS:N	8:L:207:HOH:O	1.82	1.09
2:L:4:GLY:C	8:L:216:HOH:O	1.91	1.08
1:A:8:THR:CG2	2:B:139:GLU:HA	1.82	1.08
1:K:56:LEU:CA	8:K:913:HOH:O	2.02	1.07
2:L:19:ASP:CA	8:L:228:HOH:O	1.89	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:98:LEU:N	8:B:213:HOH:O	1.83	1.07
1:A:305:THR:HG21	1:A:309:CYS:SG	1.93	1.07
2:J:124:SER:HB3	8:J:213:HOH:O	1.54	1.07
1:C:76:LEU:O	1:C:77:SER:HB2	1.54	1.06
1:C:121:GLU:OE2	1:C:258:ARG:NH1	1.88	1.06
2:H:66:VAL:HA	8:H:205:HOH:O	1.55	1.06
1:K:34:HIS:CE1	8:K:908:HOH:O	2.09	1.05
2:H:66:VAL:CA	8:H:205:HOH:O	2.04	1.05
1:K:213:SER:O	8:K:910:HOH:O	1.73	1.05
1:C:258:ARG:N	8:C:717:HOH:O	1.88	1.05
1:K:14:HIS:NE2	1:K:15:ALA:O	1.90	1.05
1:C:327:ILE:O	1:C:327:ILE:HG22	1.55	1.05
1:G:324:LEU:N	8:G:717:HOH:O	1.86	1.05
1:I:211:ARG:NH2	8:I:923:HOH:O	1.90	1.04
1:K:56:LEU:N	8:K:913:HOH:O	1.91	1.04
1:A:234:ASN:N	8:A:733:HOH:O	1.89	1.03
1:A:167:LEU:C	1:A:167:LEU:HD12	1.77	1.03
1:I:327:ILE:O	1:I:327:ILE:HG22	1.55	1.03
2:L:4:GLY:HA2	8:L:216:HOH:O	1.58	1.02
2:L:4:GLY:CA	8:L:216:HOH:O	2.08	1.02
1:K:154:LEU:HA	8:K:902:HOH:O	1.57	1.02
2:B:71:ASN:ND2	2:B:71:ASN:H	1.55	1.02
2:B:71:ASN:ND2	2:B:74:GLU:OE1	1.93	1.01
1:A:233:MET:CA	8:A:733:HOH:O	2.07	1.01
2:J:142:HIS:CE1	2:J:161:LYS:HD3	1.95	1.01
1:K:292:ILE:HG22	1:K:292:ILE:O	1.56	1.01
1:G:90:SER:CA	8:G:719:HOH:O	2.06	1.01
1:C:51:ARG:N	8:C:715:HOH:O	1.69	1.01
2:H:75:LYS:HD2	8:H:204:HOH:O	1.60	1.01
1:K:14:HIS:CD2	1:K:15:ALA:H	1.79	1.01
1:A:8:THR:HG23	2:B:139:GLU:HA	1.44	1.00
1:A:30:VAL:N	8:A:719:HOH:O	1.93	1.00
1:E:119:ARG:HG2	1:E:119:ARG:O	1.60	0.99
1:G:324:LEU:CA	8:G:717:HOH:O	2.06	0.99
1:K:33:THR:HB	8:K:908:HOH:O	1.60	0.99
1:I:190:THR:O	1:I:190:THR:HG23	1.59	0.98
1:K:14:HIS:CD2	1:K:15:ALA:N	2.30	0.98
1:A:118:GLU:OE2	8:A:720:HOH:O	1.81	0.98
1:K:13:TYR:N	8:K:925:HOH:O	1.89	0.97
1:G:324:LEU:C	8:G:717:HOH:O	2.02	0.97
1:C:245:LYS:NZ	1:I:264:GLU:OE2	1.97	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:74:GLU:O	8:K:905:HOH:O	1.81	0.97
2:B:71:ASN:N	2:B:71:ASN:HD22	1.59	0.97
1:A:91:SER:OG	8:A:732:HOH:O	1.83	0.97
1:G:317:LYS:O	8:G:710:HOH:O	1.82	0.97
1:I:190:THR:HG22	1:I:193:ASP:HB2	1.43	0.96
1:E:51:ARG:CD	8:E:718:HOH:O	2.11	0.96
1:E:294:THR:O	8:E:725:HOH:O	1.83	0.95
1:A:133:ASP:OD2	1:A:136:LYS:HD3	1.64	0.95
2:B:97:GLU:C	8:B:213:HOH:O	1.97	0.95
1:C:258:ARG:CB	8:C:717:HOH:O	1.99	0.95
1:A:102:ILE:HD12	1:A:102:ILE:N	1.79	0.95
1:C:159:LYS:HZ2	1:C:199:GLN:HE21	1.12	0.95
2:L:52:VAL:O	8:L:214:HOH:O	1.83	0.95
1:I:305:THR:H	2:J:66:VAL:HG13	1.30	0.94
1:A:12:GLY:N	2:B:14:TRP:CH2	2.36	0.94
1:A:185:ILE:O	8:A:733:HOH:O	1.84	0.94
2:J:106:ARG:HD3	8:J:214:HOH:O	1.65	0.94
2:H:103:GLU:O	8:H:211:HOH:O	1.86	0.94
1:A:91:SER:CB	8:A:717:HOH:O	2.15	0.94
1:E:66:TRP:HE1	1:E:77:SER:CB	1.80	0.93
1:K:8:THR:OG1	2:L:27:GLN:HB3	1.69	0.93
1:I:44:ASN:H	1:I:44:ASN:HD22	1.17	0.93
2:H:154:ASN:OD1	8:H:203:HOH:O	1.85	0.93
1:A:89:PRO:O	8:A:711:HOH:O	1.84	0.93
1:I:190:THR:HG23	1:I:193:ASP:H	0.76	0.92
1:E:51:ARG:HG2	8:E:718:HOH:O	1.69	0.92
1:E:66:TRP:NE1	1:E:77:SER:HB2	1.82	0.92
2:J:162:TYR:CD2	8:J:216:HOH:O	2.22	0.92
1:C:153:ASN:O	8:C:717:HOH:O	1.86	0.92
1:G:7:ASP:N	2:H:140:PHE:N	2.17	0.92
1:K:14:HIS:CG	1:K:15:ALA:N	2.29	0.92
1:C:74:GLU:O	1:C:74:GLU:HG2	1.69	0.92
2:J:106:ARG:HB2	8:J:206:HOH:O	1.67	0.91
1:A:105:GLU:O	8:A:734:HOH:O	1.88	0.91
2:J:26:HIS:HD2	2:J:26:HIS:O	1.52	0.91
1:A:29:ASN:N	8:A:719:HOH:O	2.03	0.91
1:A:133:ASP:OD2	1:A:136:LYS:CE	2.19	0.91
1:G:90:SER:N	8:G:719:HOH:O	2.04	0.90
2:L:138:PHE:O	8:L:231:HOH:O	1.89	0.90
2:J:142:HIS:CE1	2:J:161:LYS:HE2	2.07	0.90
1:E:305:THR:CB	8:E:719:HOH:O	2.20	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:19:ASP:O	8:L:228:HOH:O	1.89	0.90
1:A:29:ASN:ND2	3:A:602:NAG:O7	2.04	0.90
2:L:30:GLN:NE2	8:L:212:HOH:O	2.05	0.90
1:I:120:PHE:HB3	8:I:924:HOH:O	1.71	0.90
1:E:282:ASN:HD22	1:E:283:THR:N	1.69	0.90
2:J:103:GLU:OE1	8:J:214:HOH:O	1.89	0.90
1:K:15:ALA:HB1	8:K:904:HOH:O	1.71	0.90
1:A:102:ILE:CD1	1:A:102:ILE:N	2.30	0.90
1:K:266:ASN:ND2	1:K:266:ASN:H	1.68	0.89
2:D:161:LYS:O	8:D:204:HOH:O	1.90	0.89
1:I:190:THR:CG2	1:I:193:ASP:HB2	2.01	0.89
1:K:106:GLU:OE1	8:K:915:HOH:O	1.89	0.89
1:A:293:ASN:ND2	8:A:708:HOH:O	1.82	0.89
1:E:51:ARG:NE	8:E:718:HOH:O	2.05	0.89
1:A:102:ILE:CD1	1:A:102:ILE:H	1.85	0.89
1:A:233:MET:HA	8:A:733:HOH:O	1.66	0.89
1:I:190:THR:CG2	1:I:193:ASP:CB	2.50	0.88
1:A:233:MET:C	8:A:733:HOH:O	2.07	0.88
1:A:133:ASP:OD2	1:A:136:LYS:CD	2.21	0.88
1:C:71:PRO:HA	1:C:152:LYS:NZ	1.86	0.88
1:G:324:LEU:O	8:G:717:HOH:O	1.92	0.88
1:G:307:GLY:O	8:G:703:HOH:O	1.92	0.88
1:A:28:LYS:C	8:A:719:HOH:O	2.12	0.88
1:G:96:CYS:O	1:G:227:ARG:HD2	1.74	0.87
1:G:133:ASP:OD2	1:G:136:LYS:HG3	1.75	0.87
2:L:37:ASP:C	8:L:207:HOH:O	2.13	0.87
2:B:91:ILE:O	2:B:95:ASN:OD1	1.92	0.87
2:D:150:GLU:O	2:D:154:ASN:HB2	1.74	0.87
2:L:8:GLY:HA3	8:L:216:HOH:O	1.74	0.87
1:A:143:PRO:HD2	3:A:603:NAG:H62	1.56	0.87
1:C:11:ILE:CD1	2:D:24:TYR:HE2	1.82	0.87
1:A:308:LYS:NZ	8:A:723:HOH:O	2.06	0.87
1:C:62:ASN:OD1	8:C:729:HOH:O	1.92	0.86
1:K:66:TRP:HE1	1:K:77:SER:HB2	1.39	0.86
1:A:241:GLU:HG2	8:A:715:HOH:O	1.75	0.86
1:A:167:LEU:O	1:A:167:LEU:HD12	1.75	0.86
1:E:305:THR:OG1	8:E:719:HOH:O	1.94	0.85
1:A:185:ILE:N	8:A:721:HOH:O	2.08	0.85
1:C:204:TYR:HD2	8:C:711:HOH:O	1.59	0.85
2:L:4:GLY:O	8:L:216:HOH:O	1.90	0.85
1:I:44:ASN:HD21	1:I:291:ALA:HB3	1.42	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:265:ARG:O	8:K:909:HOH:O	1.94	0.85
1:G:114:VAL:HG11	1:G:117:PHE:HB2	1.57	0.84
1:A:133:ASP:OD2	1:A:136:LYS:HE2	1.77	0.84
1:C:114:VAL:HG11	1:C:117:PHE:HB2	1.58	0.84
1:A:114:VAL:HG11	1:A:117:PHE:HB2	1.59	0.84
1:I:172:ILE:HG23	1:I:172:ILE:O	1.78	0.84
1:G:293:ASN:ND2	8:G:706:HOH:O	2.09	0.84
1:C:327:ILE:N	1:C:327:ILE:HD12	1.92	0.84
1:K:260:ALA:N	8:K:914:HOH:O	2.10	0.84
1:I:114:VAL:HG11	1:I:117:PHE:HB2	1.57	0.84
1:G:315:SER:HA	8:H:207:HOH:O	1.78	0.84
1:I:145:ALA:O	8:I:903:HOH:O	1.96	0.84
1:I:194:GLN:OE1	1:I:200:ASN:O	1.96	0.84
2:L:150:GLU:O	2:L:154:ASN:HB2	1.78	0.84
1:I:114:VAL:CG1	1:I:116:SER:O	2.27	0.83
1:I:118:GLU:OE1	8:I:913:HOH:O	1.95	0.83
2:L:53:ASN:HA	8:L:214:HOH:O	1.77	0.83
2:H:127:LYS:HD2	2:H:127:LYS:H	1.43	0.83
1:A:159:LYS:NZ	1:A:195:GLN:O	2.11	0.83
1:C:11:ILE:HD12	2:D:24:TYR:CE2	2.12	0.83
2:F:136:GLY:N	8:F:201:HOH:O	2.11	0.83
2:H:150:GLU:O	2:H:154:ASN:HB2	1.79	0.83
2:H:1:GLY:HA2	8:H:209:HOH:O	1.77	0.83
1:I:144:HIS:O	8:I:903:HOH:O	1.96	0.83
2:J:132:GLU:OE1	8:J:212:HOH:O	1.95	0.83
8:H:213:HOH:O	2:J:66:VAL:HG23	1.78	0.82
2:J:150:GLU:O	2:J:154:ASN:HB2	1.79	0.82
2:F:127:LYS:H	2:F:127:LYS:HD2	1.44	0.82
1:C:59:GLY:N	8:C:712:HOH:O	2.11	0.82
1:I:76:LEU:O	1:I:77:SER:HB2	1.79	0.82
2:J:127:LYS:H	2:J:127:LYS:HD2	1.42	0.82
1:K:74:GLU:C	8:K:905:HOH:O	2.13	0.82
1:I:74:GLU:OE1	8:I:914:HOH:O	1.97	0.82
2:F:150:GLU:O	2:F:154:ASN:HB2	1.79	0.82
2:J:123:ARG:NH1	8:J:213:HOH:O	2.11	0.82
1:A:326:ASN:C	1:A:327:ILE:CG1	2.46	0.82
2:B:127:LYS:HD2	2:B:127:LYS:H	1.42	0.82
1:K:266:ASN:HD22	1:K:266:ASN:H	1.24	0.82
2:L:19:ASP:N	8:L:226:HOH:O	2.10	0.81
2:B:150:GLU:O	2:B:154:ASN:HB2	1.79	0.81
1:C:305:THR:H	2:D:66:VAL:HG13	1.46	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:231:GLY:O	1:E:232:ARG:HD3	1.80	0.81
2:J:26:HIS:C	2:J:26:HIS:CD2	2.54	0.81
1:E:13:TYR:CE2	2:F:6:ILE:HA	2.16	0.81
1:K:259:TYR:CZ	8:K:906:HOH:O	2.33	0.81
1:K:114:VAL:HG21	1:K:117:PHE:HB2	1.63	0.80
2:H:120:GLU:OE1	2:H:123:ARG:CZ	2.30	0.80
1:C:62:ASN:CG	8:C:729:HOH:O	2.20	0.80
2:D:127:LYS:H	2:D:127:LYS:HD2	1.43	0.80
1:K:11:ILE:O	1:K:11:ILE:HG22	1.81	0.80
1:K:258:ARG:N	8:K:902:HOH:O	2.12	0.80
1:C:69:GLY:O	1:C:152:LYS:HG2	1.82	0.80
2:H:106:ARG:HB2	8:H:211:HOH:O	1.80	0.80
1:C:195:GLN:HA	1:C:198:TYR:O	1.82	0.80
1:A:132:HIS:HD2	8:A:729:HOH:O	1.64	0.80
1:I:190:THR:HG21	1:I:193:ASP:CG	2.02	0.80
2:D:124:SER:O	2:D:127:LYS:HE3	1.81	0.80
1:K:51:ARG:NH1	1:K:51:ARG:HB3	1.96	0.80
1:I:190:THR:O	1:I:190:THR:CG2	2.30	0.80
1:G:41:ASP:O	8:G:708:HOH:O	1.98	0.80
1:I:172:ILE:CG2	1:I:172:ILE:O	2.30	0.80
1:E:114:VAL:HG21	1:E:117:PHE:HB2	1.63	0.80
1:K:7:ASP:O	8:K:911:HOH:O	2.00	0.79
1:C:172:ILE:O	8:C:722:HOH:O	2.00	0.79
2:L:148:CYS:O	8:L:223:HOH:O	1.99	0.79
1:E:244:ASP:OD1	8:E:728:HOH:O	1.99	0.79
1:C:326:ASN:C	1:C:327:ILE:HD12	2.03	0.79
1:A:195:GLN:HG2	8:A:704:HOH:O	1.82	0.79
1:E:319:ARG:CG	1:E:319:ARG:HH11	1.94	0.79
1:A:167:LEU:C	1:A:167:LEU:CD1	2.50	0.79
1:C:159:LYS:NZ	1:C:199:GLN:HE21	1.79	0.79
1:E:11:ILE:HD12	2:F:119:TYR:HA	1.62	0.79
2:F:123:ARG:NH2	8:F:204:HOH:O	1.96	0.79
1:C:100:ASP:OD1	8:C:720:HOH:O	2.00	0.79
1:E:110:GLN:OE1	8:E:731:HOH:O	2.01	0.79
1:E:72:GLU:HG2	6:E:601:NAG:H82	1.62	0.79
1:G:7:ASP:N	1:G:8:THR:HA	1.98	0.78
1:C:50:LEU:CA	8:C:715:HOH:O	2.32	0.78
4:A:607:NAG:O7	8:A:706:HOH:O	2.01	0.78
2:B:14:TRP:HE3	2:B:17:MET:HE2	1.48	0.78
1:K:266:ASN:ND2	1:K:266:ASN:N	2.30	0.78
1:I:190:THR:CG2	1:I:193:ASP:CG	2.51	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:ILE:HG12	8:A:721:HOH:O	1.82	0.78
1:A:61:CYS:O	8:A:717:HOH:O	2.00	0.78
1:G:143:PRO:HD2	3:G:602:NAG:H62	1.66	0.78
1:E:203:THR:O	1:E:217:LYS:CE	2.31	0.78
1:K:13:TYR:O	8:K:925:HOH:O	2.01	0.77
2:J:142:HIS:CE1	2:J:161:LYS:CD	2.67	0.77
2:L:127:LYS:H	2:L:127:LYS:HD2	1.47	0.77
1:C:76:LEU:O	1:C:77:SER:CB	2.29	0.77
1:I:244:ASP:OD1	8:I:925:HOH:O	2.01	0.77
1:E:282:ASN:OD1	8:E:710:HOH:O	2.03	0.77
1:E:11:ILE:CD1	2:F:119:TYR:HA	2.15	0.77
1:A:185:ILE:CB	8:A:721:HOH:O	2.33	0.77
2:D:109:ASP:OD2	8:F:205:HOH:O	2.02	0.77
2:J:18:VAL:HB	8:J:211:HOH:O	1.85	0.77
2:D:150:GLU:OE2	2:D:150:GLU:CA	2.30	0.77
1:C:59:GLY:O	8:C:712:HOH:O	2.01	0.77
1:K:11:ILE:O	1:K:11:ILE:CG2	2.33	0.77
2:J:18:VAL:O	8:J:211:HOH:O	2.02	0.77
1:E:47:LEU:HD12	1:E:286:GLN:NE2	2.00	0.77
2:J:26:HIS:O	2:J:26:HIS:CD2	2.36	0.76
1:K:294:THR:N	8:K:916:HOH:O	1.82	0.76
1:I:114:VAL:HG13	1:I:116:SER:O	1.86	0.76
2:H:6:ILE:O	8:H:209:HOH:O	2.01	0.76
1:E:319:ARG:HH11	1:E:319:ARG:HG2	1.49	0.76
1:A:107:LEU:CD2	1:A:107:LEU:C	2.53	0.76
1:I:89:PRO:HD2	8:I:908:HOH:O	1.85	0.76
1:E:51:ARG:CG	8:E:718:HOH:O	2.23	0.76
1:A:192:ALA:O	1:A:196:SER:OG	2.03	0.76
1:K:27:GLU:OE2	8:K:921:HOH:O	2.04	0.76
1:K:56:LEU:HB2	8:K:913:HOH:O	1.85	0.76
1:C:64:ALA:N	8:C:729:HOH:O	2.12	0.75
1:I:105:GLU:O	8:I:906:HOH:O	2.02	0.75
1:K:246:ILE:HG13	1:K:247:THR:N	2.00	0.75
1:I:190:THR:HG22	1:I:193:ASP:CB	2.13	0.75
1:K:7:ASP:HA	2:L:27:GLN:O	1.86	0.75
1:A:159:LYS:NZ	1:A:199:GLN:NE2	2.33	0.75
1:A:130:PRO:HD2	8:A:729:HOH:O	1.87	0.75
1:K:305:THR:H	2:L:66:VAL:HG13	1.50	0.75
1:I:120:PHE:CA	8:I:924:HOH:O	2.33	0.75
1:I:241:GLU:HG2	8:I:925:HOH:O	1.87	0.75
1:K:24:THR:HG22	2:L:104:ASN:HB3	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:11:ILE:HD13	2:D:24:TYR:HD2	1.00	0.74
2:J:106:ARG:CB	8:J:206:HOH:O	2.29	0.74
2:H:107:THR:N	8:H:211:HOH:O	2.20	0.74
1:A:8:THR:HG21	2:B:138:PHE:O	1.87	0.74
1:C:222:ILE:HG23	1:C:230:GLU:HG2	1.69	0.74
1:I:119:ARG:NH1	8:I:919:HOH:O	2.21	0.74
1:A:76:LEU:O	8:A:739:HOH:O	2.05	0.74
1:I:282:ASN:ND2	8:I:902:HOH:O	2.18	0.74
1:A:7:ASP:C	2:B:27:GLN:O	2.26	0.73
1:A:102:ILE:HD13	1:A:102:ILE:H	1.51	0.73
1:G:315:SER:CA	8:H:207:HOH:O	2.34	0.73
2:H:162:TYR:OH	8:H:206:HOH:O	2.05	0.73
1:C:71:PRO:HA	1:C:152:LYS:HZ2	1.53	0.73
1:A:82:TRP:O	8:A:740:HOH:O	2.05	0.73
1:G:90:SER:HA	8:G:719:HOH:O	1.76	0.73
1:I:327:ILE:O	1:I:327:ILE:CG2	2.30	0.73
1:E:8:THR:OG1	8:E:712:HOH:O	2.06	0.73
1:C:119:ARG:HB2	1:C:261:PHE:CD2	2.24	0.72
2:B:4:GLY:CA	8:B:212:HOH:O	2.37	0.72
2:J:142:HIS:CE1	2:J:161:LYS:CE	2.72	0.72
2:H:83:LYS:NZ	8:H:213:HOH:O	1.91	0.72
1:I:189:SER:HB2	1:I:222:ILE:HD13	1.71	0.72
1:K:294:THR:O	8:K:916:HOH:O	2.08	0.72
2:F:106:ARG:NH2	8:F:205:HOH:O	1.86	0.72
2:B:123:ARG:HD2	2:B:132:GLU:OE1	1.89	0.72
1:A:305:THR:HG21	1:A:309:CYS:CB	2.19	0.72
1:K:55:PRO:C	8:K:913:HOH:O	2.19	0.72
1:G:323:GLY:C	8:G:717:HOH:O	2.24	0.72
2:J:103:GLU:HA	8:J:206:HOH:O	1.90	0.72
1:A:24:THR:HG22	2:B:104:ASN:HB3	1.72	0.72
1:I:275:ASP:OD2	8:I:908:HOH:O	2.07	0.72
1:C:46:LYS:NZ	8:C:719:HOH:O	2.23	0.72
1:C:71:PRO:HA	1:C:152:LYS:HZ1	1.54	0.71
2:D:150:GLU:HA	2:D:150:GLU:OE2	1.90	0.71
1:I:118:GLU:CD	8:I:913:HOH:O	2.29	0.71
1:I:120:PHE:CB	8:I:924:HOH:O	2.32	0.71
1:E:10:CYS:HA	2:F:137:CYS:HA	1.72	0.71
2:D:132:GLU:HG2	2:D:138:PHE:HE2	1.55	0.71
1:C:130:PRO:O	1:C:160:LYS:NZ	2.23	0.71
2:B:14:TRP:HE3	2:B:17:MET:CE	2.04	0.71
2:L:132:GLU:HG2	2:L:138:PHE:HE2	1.54	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:12:GLY:O	8:B:202:HOH:O	2.07	0.71
2:J:142:HIS:HE1	2:J:161:LYS:HE2	1.54	0.71
2:J:103:GLU:O	8:J:206:HOH:O	2.08	0.71
2:B:132:GLU:HG2	2:B:138:PHE:HE2	1.56	0.70
2:B:22:TYR:OH	2:B:111:HIS:ND1	2.20	0.70
2:J:132:GLU:HG2	2:J:138:PHE:HE2	1.55	0.70
1:A:11:ILE:C	2:B:14:TRP:CH2	2.64	0.70
2:F:132:GLU:HG2	2:F:138:PHE:HE2	1.55	0.70
1:K:10:CYS:O	2:L:24:TYR:HA	1.91	0.70
2:L:157:TYR:CD1	8:L:223:HOH:O	2.44	0.70
2:D:71:ASN:OD1	2:D:71:ASN:C	2.29	0.70
2:H:132:GLU:HG2	2:H:138:PHE:HE2	1.56	0.70
2:B:71:ASN:H	2:B:71:ASN:HD22	0.79	0.70
1:A:196:SER:HB2	4:A:607:NAG:H82	1.73	0.70
1:K:292:ILE:HD11	1:K:301:ILE:HD12	1.73	0.70
1:A:145:ALA:C	1:A:147:ALA:H	1.95	0.70
1:G:11:ILE:HD12	2:H:119:TYR:HA	1.73	0.70
1:K:51:ARG:HH11	1:K:51:ARG:HB3	1.57	0.70
1:E:282:ASN:HD22	1:E:283:THR:H	1.40	0.70
1:A:191:SER:O	8:A:704:HOH:O	2.09	0.69
2:J:18:VAL:CA	8:J:211:HOH:O	2.40	0.69
1:A:313:VAL:HG12	1:A:315:SER:H	1.57	0.69
1:C:327:ILE:O	1:C:327:ILE:CG2	2.30	0.69
1:A:30:VAL:HG23	8:A:719:HOH:O	1.92	0.69
1:A:27:GLU:OE1	1:A:325:ARG:NH2	2.25	0.69
1:A:236:TYR:OH	8:A:726:HOH:O	2.08	0.69
1:G:7:ASP:N	2:H:140:PHE:HB2	2.07	0.69
1:G:89:PRO:C	8:G:719:HOH:O	2.29	0.69
1:G:313:VAL:HG12	1:G:315:SER:H	1.58	0.69
1:A:202:ASP:OD1	1:A:202:ASP:C	2.30	0.69
1:I:101:PHE:O	8:I:920:HOH:O	2.11	0.69
1:E:282:ASN:ND2	1:E:283:THR:N	2.40	0.69
2:L:139:GLU:HB3	8:L:225:HOH:O	1.91	0.69
1:A:204:TYR:CD1	1:A:204:TYR:C	2.65	0.69
1:A:74:GLU:HA	8:A:737:HOH:O	1.92	0.69
1:A:233:MET:CG	8:A:733:HOH:O	2.41	0.68
1:K:313:VAL:HG12	1:K:315:SER:H	1.58	0.68
1:C:102:ILE:HG13	1:C:236:TYR:CE2	2.28	0.68
1:E:313:VAL:HG12	1:E:315:SER:H	1.59	0.68
1:G:102:ILE:HG13	1:G:236:TYR:CE2	2.29	0.68
1:K:259:TYR:OH	8:K:906:HOH:O	2.09	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:50:LEU:CB	8:C:715:HOH:O	2.41	0.68
1:E:119:ARG:CG	1:E:119:ARG:O	2.32	0.68
2:L:97:GLU:OE2	8:L:220:HOH:O	2.12	0.68
1:K:102:ILE:HG13	1:K:236:TYR:CE2	2.29	0.68
1:I:15:ALA:O	8:I:909:HOH:O	2.11	0.68
1:I:313:VAL:HG12	1:I:315:SER:H	1.58	0.68
1:A:49:LYS:O	1:A:283:THR:HG22	1.94	0.68
1:A:120:PHE:HB3	8:A:720:HOH:O	1.93	0.68
1:E:245:LYS:NZ	8:E:721:HOH:O	2.22	0.67
1:C:11:ILE:CD1	2:D:24:TYR:HD2	1.83	0.67
1:I:190:THR:HG22	1:I:193:ASP:OD1	1.93	0.67
8:E:712:HOH:O	2:F:27:GLN:HB3	1.95	0.67
1:K:324:LEU:HD21	2:L:21:TRP:CD1	2.30	0.67
2:H:42:GLN:OE1	8:H:201:HOH:O	2.12	0.67
1:A:305:THR:HG23	1:A:309:CYS:SG	2.31	0.67
1:A:185:ILE:HB	8:A:721:HOH:O	1.92	0.67
2:L:38:LEU:N	8:L:207:HOH:O	2.25	0.67
1:E:119:ARG:HB2	1:E:261:PHE:CD2	2.30	0.67
1:I:172:ILE:HD12	1:I:245:LYS:HB3	1.76	0.67
1:A:132:HIS:CD2	8:A:729:HOH:O	2.43	0.67
1:C:313:VAL:HG12	1:C:315:SER:H	1.59	0.67
1:G:13:TYR:CE2	2:H:6:ILE:HA	2.28	0.67
1:E:102:ILE:HG13	1:E:236:TYR:CE2	2.28	0.67
1:A:128:SER:O	1:A:130:PRO:HD3	1.94	0.67
1:G:268:GLY:HA2	8:G:709:HOH:O	1.95	0.67
1:C:223:ARG:O	1:C:230:GLU:HG3	1.95	0.66
2:H:51:LYS:NZ	8:H:211:HOH:O	2.26	0.66
1:A:107:LEU:C	1:A:107:LEU:HD22	2.15	0.66
1:A:217:LYS:O	8:A:735:HOH:O	2.13	0.66
1:A:105:GLU:HA	1:A:108:ARG:HD3	1.78	0.66
1:C:49:LYS:O	1:C:283:THR:HG22	1.96	0.66
2:F:83:LYS:O	8:F:207:HOH:O	2.13	0.66
1:E:178:GLU:CD	1:E:265:ARG:NH1	2.48	0.66
2:L:37:ASP:O	8:L:207:HOH:O	2.11	0.66
1:I:194:GLN:HG3	1:I:195:GLN:N	2.10	0.66
1:I:311:LYS:HG3	2:J:92:TRP:CE2	2.31	0.66
1:A:102:ILE:HD13	1:A:235:TYR:O	1.96	0.66
2:B:71:ASN:HD21	2:B:74:GLU:CD	1.99	0.66
1:K:49:LYS:O	1:K:283:THR:HG22	1.95	0.66
1:A:76:LEU:N	8:A:738:HOH:O	1.80	0.66
1:I:102:ILE:HG13	1:I:236:TYR:CE2	2.30	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:THR:HG23	1:A:324:LEU:O	1.96	0.65
1:A:282:ASN:OD1	8:A:724:HOH:O	2.13	0.65
1:I:44:ASN:N	1:I:44:ASN:HD22	1.89	0.65
1:G:109:GLU:OE2	2:L:79:ASN:ND2	2.29	0.65
1:I:307:GLY:HA2	2:J:63:PHE:CE1	2.31	0.65
1:I:190:THR:CG2	1:I:193:ASP:N	2.38	0.65
2:J:162:TYR:CE2	8:J:216:HOH:O	2.43	0.65
1:I:265:ARG:NE	8:I:910:HOH:O	2.28	0.65
1:C:24:THR:HG22	2:D:104:ASN:HB3	1.78	0.65
1:I:25:VAL:HG12	2:L:51:LYS:HA	1.77	0.65
1:I:297:PRO:HG3	2:J:56:ILE:HA	1.78	0.65
1:I:49:LYS:O	1:I:283:THR:HG22	1.96	0.65
2:J:103:GLU:C	8:J:206:HOH:O	2.35	0.65
2:B:51:LYS:HG3	1:E:25:VAL:HG12	1.78	0.65
1:I:145:ALA:CB	8:I:903:HOH:O	2.16	0.65
1:E:178:GLU:CD	1:E:265:ARG:HH12	1.98	0.65
1:C:50:LEU:N	8:C:715:HOH:O	2.30	0.65
1:C:25:VAL:CG2	2:D:102:LEU:HD12	2.26	0.65
1:K:43:HIS:HB3	1:K:301:ILE:HD13	1.79	0.65
1:I:44:ASN:HD21	1:I:291:ALA:CB	2.09	0.65
1:E:119:ARG:HB2	1:E:261:PHE:CE2	2.32	0.65
1:A:307:GLY:HA2	2:B:63:PHE:CE1	2.32	0.65
1:C:155:ILE:HD11	1:C:258:ARG:HD3	1.80	0.64
1:G:11:ILE:CD1	2:H:119:TYR:HA	2.27	0.64
1:A:13:TYR:HD1	8:A:730:HOH:O	1.80	0.64
1:G:49:LYS:O	1:G:283:THR:HG22	1.98	0.64
2:L:139:GLU:N	8:L:225:HOH:O	2.21	0.64
1:K:16:ASN:ND2	8:K:908:HOH:O	2.05	0.64
1:E:294:THR:O	8:E:704:HOH:O	2.15	0.64
1:C:74:GLU:O	1:C:74:GLU:CG	2.28	0.64
1:C:231:GLY:HA3	8:C:701:HOH:O	1.96	0.64
1:K:293:ASN:O	1:K:293:ASN:ND2	2.30	0.64
2:J:159:TYR:N	2:J:160:PRO:CD	2.61	0.64
2:L:157:TYR:CE1	8:L:223:HOH:O	2.50	0.64
1:E:319:ARG:CG	1:E:319:ARG:NH1	2.55	0.64
1:A:107:LEU:O	1:A:107:LEU:CD2	2.46	0.64
1:C:131:ASN:HD21	1:I:243:GLY:HA3	1.62	0.64
2:F:19:ASP:N	2:F:19:ASP:OD1	2.30	0.64
1:E:50:LEU:HD21	1:E:306:ILE:HG22	1.80	0.64
1:K:82:TRP:HB3	8:K:913:HOH:O	1.97	0.64
1:I:25:VAL:CG1	2:L:51:LYS:HG3	2.28	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:131:LYS:O	8:L:225:HOH:O	2.16	0.64
1:E:257:PRO:HB3	8:E:726:HOH:O	1.98	0.64
1:C:204:TYR:CD2	8:C:711:HOH:O	2.41	0.63
1:A:108:ARG:NH1	8:A:709:HOH:O	2.10	0.63
1:I:14:HIS:HE1	2:J:18:VAL:HA	1.64	0.63
2:L:123:ARG:C	8:L:205:HOH:O	2.36	0.63
1:K:66:TRP:NE1	1:K:77:SER:HB2	2.12	0.63
1:C:119:ARG:HB2	1:C:261:PHE:CE2	2.33	0.63
1:C:59:GLY:CA	8:C:712:HOH:O	2.45	0.63
2:J:18:VAL:CB	8:J:211:HOH:O	2.45	0.63
1:C:133:ASP:O	8:C:728:HOH:O	2.15	0.63
1:K:66:TRP:HE1	1:K:77:SER:CB	2.09	0.63
1:C:327:ILE:N	1:C:327:ILE:CD1	2.60	0.63
2:B:14:TRP:CE3	2:B:17:MET:CE	2.82	0.63
1:E:43:HIS:HB3	1:E:301:ILE:HD13	1.80	0.62
1:C:194:GLN:O	1:C:198:TYR:O	2.17	0.62
1:C:64:ALA:HB3	8:C:729:HOH:O	1.98	0.62
1:I:43:HIS:HB3	1:I:301:ILE:HD13	1.80	0.62
1:A:43:HIS:HB3	1:A:301:ILE:HD13	1.81	0.62
1:A:224:PRO:HG2	1:C:209:SER:HA	1.79	0.62
1:K:106:GLU:O	1:K:110:GLN:HG2	2.00	0.62
1:A:159:LYS:HZ2	1:A:199:GLN:NE2	1.95	0.62
2:L:131:LYS:HG3	8:L:225:HOH:O	2.00	0.62
1:I:231:GLY:O	1:I:232:ARG:NH1	2.32	0.62
1:A:305:THR:HG22	1:A:309:CYS:SG	2.37	0.62
1:E:13:TYR:CD2	2:F:6:ILE:HG12	2.35	0.62
1:A:131:ASN:O	1:A:132:HIS:CG	2.53	0.62
1:A:145:ALA:O	1:A:147:ALA:N	2.32	0.62
1:I:106:GLU:O	1:I:110:GLN:HG2	1.99	0.62
1:G:43:HIS:HB3	1:G:301:ILE:HD13	1.82	0.62
1:E:229:GLN:NE2	5:E:603:SIA:O1A	2.30	0.62
2:L:158:ASP:O	2:L:160:PRO:HD3	2.00	0.62
1:G:32:VAL:HB	8:G:717:HOH:O	2.00	0.62
1:K:25:VAL:CG2	2:L:102:LEU:HD12	2.29	0.62
1:G:106:GLU:O	1:G:110:GLN:HG2	2.00	0.62
1:C:43:HIS:HB3	1:C:301:ILE:HD13	1.80	0.61
2:B:85:ASP:OD1	2:F:83:LYS:NZ	2.33	0.61
1:G:10:CYS:HA	2:H:137:CYS:HA	1.82	0.61
1:K:8:THR:HG1	2:L:27:GLN:HB3	1.65	0.61
1:E:49:LYS:O	1:E:283:THR:HG22	2.01	0.61
1:G:297:PRO:HG3	2:H:56:ILE:HA	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:THR:CG2	2:B:139:GLU:CA	2.72	0.61
1:K:239:LEU:CD1	1:K:265:ARG:NH1	2.63	0.61
1:I:141:ALA:O	1:I:143:PRO:HD3	2.00	0.61
2:L:139:GLU:CB	8:L:225:HOH:O	2.47	0.61
1:K:105:GLU:OE2	2:L:71:ASN:HB3	1.99	0.61
1:C:11:ILE:HD11	2:D:24:TYR:CE2	2.31	0.61
1:G:195:GLN:HA	1:G:198:TYR:O	2.00	0.61
2:B:62:GLN:OE1	1:E:314:LYS:NZ	2.20	0.61
2:D:159:TYR:N	2:D:160:PRO:CD	2.64	0.61
1:G:133:ASP:OD2	1:G:136:LYS:CG	2.48	0.61
2:B:79:ASN:ND2	1:C:109:GLU:OE2	2.34	0.61
3:A:602:NAG:C1	3:A:602:NAG:O7	2.49	0.60
1:I:50:LEU:HD11	1:I:306:ILE:HG22	1.83	0.60
1:I:145:ALA:C	8:I:903:HOH:O	2.36	0.60
1:A:77:SER:N	8:A:738:HOH:O	2.30	0.60
1:G:161:GLY:O	1:G:162:ASN:HB2	2.01	0.60
1:E:161:GLY:O	1:E:162:ASN:HB2	2.00	0.60
1:K:9:LEU:HD21	2:L:153:LYS:HE2	1.83	0.60
1:G:90:SER:HB2	8:G:719:HOH:O	2.01	0.60
2:D:19:ASP:HB2	2:D:36:ALA:HB3	1.84	0.60
2:L:152:VAL:O	8:L:227:HOH:O	2.17	0.60
2:D:162:TYR:CZ	8:D:205:HOH:O	2.51	0.60
2:B:95:ASN:HD21	2:D:95:ASN:HD22	1.49	0.60
1:A:326:ASN:O	1:A:327:ILE:CG1	2.35	0.60
2:L:124:SER:HA	8:L:205:HOH:O	2.01	0.60
1:G:141:ALA:O	1:G:143:PRO:HD3	2.01	0.60
2:D:71:ASN:ND2	8:D:207:HOH:O	2.33	0.60
1:E:25:VAL:CG2	2:F:102:LEU:HD12	2.31	0.60
1:K:76:LEU:N	8:K:905:HOH:O	2.22	0.60
1:E:141:ALA:O	1:E:143:PRO:HD3	2.02	0.60
1:G:8:THR:HG22	2:H:139:GLU:HA	1.84	0.59
2:J:161:LYS:O	8:J:207:HOH:O	2.16	0.59
1:E:211:ARG:HD2	8:E:728:HOH:O	2.00	0.59
1:I:14:HIS:CE1	2:J:18:VAL:HA	2.37	0.59
1:K:141:ALA:O	1:K:143:PRO:HD3	2.02	0.59
1:C:299:GLN:HG2	1:C:310:PRO:HB2	1.83	0.59
2:H:43:ASN:O	2:H:47:GLU:HG3	2.02	0.59
1:K:312:TYR:CD2	2:L:89:LEU:HD13	2.36	0.59
1:I:44:ASN:ND2	1:I:291:ALA:HB3	2.14	0.59
2:D:150:GLU:N	2:D:150:GLU:OE2	2.35	0.59
1:C:141:ALA:O	1:C:143:PRO:HD3	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:19:ASP:HB2	2:J:36:ALA:HB3	1.85	0.59
1:K:56:LEU:CB	8:K:913:HOH:O	2.32	0.59
2:L:23:GLY:HA3	2:L:36:ALA:HA	1.84	0.59
1:C:168:SER:HB3	1:I:176:GLY:HA2	1.85	0.59
1:C:25:VAL:HG22	2:D:102:LEU:HD12	1.84	0.59
1:E:58:LEU:HD11	1:E:63:ILE:HD13	1.84	0.59
1:E:265:ARG:HG3	8:E:720:HOH:O	2.02	0.59
1:A:214:LYS:HG2	1:A:216:PHE:CZ	2.36	0.59
1:A:131:ASN:C	1:A:132:HIS:CD2	2.77	0.59
1:E:8:THR:O	8:E:712:HOH:O	2.17	0.59
1:I:25:VAL:HG12	2:L:51:LYS:HG3	1.85	0.59
1:C:107:LEU:HD12	1:C:107:LEU:O	2.03	0.59
2:H:19:ASP:HB2	2:H:36:ALA:HB3	1.84	0.59
2:B:19:ASP:HB2	2:B:36:ALA:HB3	1.84	0.59
2:B:67:GLY:C	2:B:68:LYS:HG2	2.23	0.59
1:C:231:GLY:CA	8:C:701:HOH:O	2.51	0.58
1:A:26:LEU:HB3	2:D:47:GLU:HB3	1.85	0.58
1:I:312:TYR:CD2	2:J:89:LEU:HD13	2.38	0.58
2:B:159:TYR:N	2:B:160:PRO:HD3	2.18	0.58
1:E:281:CYS:N	8:E:709:HOH:O	2.35	0.58
1:K:13:TYR:OH	2:L:12:GLY:N	2.32	0.58
1:G:315:SER:CB	8:H:207:HOH:O	2.51	0.58
2:L:19:ASP:HB2	2:L:36:ALA:HB3	1.85	0.58
6:E:601:NAG:H61	6:E:602:NAG:C7	2.33	0.58
1:I:58:LEU:HD11	1:I:63:ILE:HD13	1.85	0.58
2:L:2:LEU:N	8:L:222:HOH:O	2.36	0.58
2:L:132:GLU:HG2	2:L:138:PHE:CE2	2.38	0.58
2:L:123:ARG:HD2	2:L:132:GLU:OE1	2.04	0.58
1:E:8:THR:HG22	2:F:139:GLU:HA	1.84	0.58
1:G:92:ASP:HB2	3:G:602:NAG:H81	1.86	0.58
1:C:162:ASN:HD22	1:C:199:GLN:NE2	2.02	0.58
1:C:25:VAL:HG12	2:F:51:LYS:HG3	1.85	0.58
1:I:161:GLY:O	1:I:162:ASN:HB2	2.02	0.58
5:C:602:SIA:H6	5:C:602:SIA:O1A	2.03	0.58
1:E:44:ASN:HD21	1:E:291:ALA:H	1.51	0.58
1:A:50:LEU:HD11	1:A:306:ILE:HG22	1.86	0.58
2:D:123:ARG:HD2	2:D:132:GLU:OE1	2.03	0.58
1:I:40:GLU:OE1	1:I:294:THR:OG1	2.19	0.58
1:C:158:VAL:HG22	8:C:728:HOH:O	2.03	0.58
1:G:50:LEU:HD11	1:G:306:ILE:HG22	1.86	0.58
1:K:58:LEU:HD11	1:K:63:ILE:HD13	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:44:ASN:HD21	1:G:291:ALA:H	1.52	0.58
1:C:11:ILE:HD12	2:D:24:TYR:HE2	1.54	0.57
1:A:185:ILE:CG1	8:A:721:HOH:O	2.42	0.57
1:A:29:ASN:C	8:A:719:HOH:O	2.34	0.57
1:I:44:ASN:ND2	1:I:44:ASN:H	1.96	0.57
1:K:245:LYS:HG3	1:K:246:ILE:N	2.18	0.57
2:B:123:ARG:HG2	2:B:123:ARG:O	2.00	0.57
1:A:15:ALA:HB2	8:A:712:HOH:O	2.04	0.57
1:A:107:LEU:C	1:A:107:LEU:HD23	2.24	0.57
2:B:51:LYS:HG3	1:E:25:VAL:CG1	2.34	0.57
1:K:312:TYR:HD2	2:L:89:LEU:HD13	1.70	0.57
1:A:107:LEU:O	1:A:107:LEU:HD23	2.05	0.57
1:A:161:GLY:O	1:A:162:ASN:HB2	2.03	0.57
1:C:107:LEU:HD12	1:C:107:LEU:C	2.24	0.57
1:K:302:HIS:CE1	1:K:304:ILE:HB	2.40	0.57
1:I:293:ASN:ND2	8:I:905:HOH:O	2.36	0.57
2:L:18:VAL:CA	8:L:226:HOH:O	2.52	0.57
1:C:153:ASN:C	8:C:717:HOH:O	2.36	0.57
2:J:159:TYR:N	2:J:160:PRO:HD2	2.20	0.57
1:K:75:SER:CA	8:K:905:HOH:O	2.52	0.57
1:C:172:ILE:CG2	8:C:722:HOH:O	2.52	0.57
1:C:106:GLU:O	1:C:110:GLN:HG2	2.05	0.57
1:A:225:LYS:NZ	4:A:606:GAL:O3	2.37	0.57
2:L:123:ARG:HD2	8:L:217:HOH:O	2.05	0.57
1:C:14:HIS:N	2:D:21:TRP:O	2.37	0.57
1:K:13:TYR:CE1	2:L:12:GLY:O	2.58	0.57
1:A:145:ALA:C	1:A:147:ALA:N	2.58	0.57
1:A:28:LYS:NZ	3:A:602:NAG:O6	2.32	0.57
1:A:91:SER:HB3	8:A:717:HOH:O	1.94	0.57
1:A:313:VAL:CG1	1:A:315:SER:H	2.18	0.57
2:D:23:GLY:HA3	2:D:36:ALA:HA	1.86	0.57
2:D:69:GLU:N	8:D:208:HOH:O	1.86	0.57
1:A:44:ASN:HD21	1:A:291:ALA:H	1.52	0.57
2:B:26:HIS:O	2:B:26:HIS:HD2	1.88	0.57
2:B:26:HIS:O	2:B:26:HIS:CD2	2.58	0.57
1:I:327:ILE:N	1:I:327:ILE:CD1	2.67	0.57
1:I:25:VAL:O	2:L:51:LYS:HA	2.05	0.57
2:J:23:GLY:HA3	2:J:36:ALA:HA	1.87	0.57
1:K:44:ASN:HD21	1:K:291:ALA:H	1.51	0.57
1:E:48:CYS:HB3	1:E:281:CYS:O	2.04	0.56
2:L:124:SER:CA	8:L:205:HOH:O	2.53	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:58:LEU:HD11	1:C:63:ILE:HD13	1.86	0.56
1:A:188:PRO:HG2	1:A:194:GLN:NE2	2.20	0.56
1:I:173:ASN:ND2	1:I:179:VAL:HG23	2.19	0.56
2:H:23:GLY:HA3	2:H:36:ALA:HA	1.87	0.56
2:L:19:ASP:C	8:L:228:HOH:O	2.11	0.56
1:K:13:TYR:CE1	2:L:12:GLY:C	2.78	0.56
1:E:305:THR:H	2:F:66:VAL:HG13	1.70	0.56
2:D:160:PRO:O	2:D:161:LYS:HG2	2.05	0.56
1:K:10:CYS:O	2:L:24:TYR:CA	2.53	0.56
1:C:297:PRO:HG3	2:D:56:ILE:HA	1.87	0.56
1:K:259:TYR:HA	8:K:914:HOH:O	2.05	0.56
1:C:155:ILE:HG22	1:C:157:LEU:CD1	2.36	0.56
1:E:11:ILE:O	2:F:10:ILE:HD13	2.05	0.56
2:F:123:ARG:HD2	2:F:132:GLU:OE1	2.05	0.56
1:K:155:ILE:HD11	1:K:258:ARG:HD2	1.87	0.56
2:B:4:GLY:N	8:B:212:HOH:O	2.37	0.56
1:K:259:TYR:CA	8:K:914:HOH:O	2.54	0.56
1:K:13:TYR:O	2:L:13:GLY:HA2	2.05	0.56
1:E:305:THR:HB	8:E:719:HOH:O	1.98	0.56
1:I:224:PRO:HG2	1:K:209:SER:HA	1.88	0.56
1:K:44:ASN:ND2	1:K:291:ALA:H	2.04	0.56
1:E:188:PRO:HG2	1:E:194:GLN:NE2	2.21	0.56
1:I:133:ASP:HB3	1:I:158:VAL:HG23	1.88	0.56
1:G:188:PRO:HG2	1:G:194:GLN:NE2	2.21	0.56
1:C:29:ASN:ND2	8:C:706:HOH:O	2.09	0.56
1:A:23:ASP:O	2:B:104:ASN:ND2	2.22	0.56
1:I:302:HIS:CE1	1:I:304:ILE:HB	2.41	0.56
1:A:302:HIS:CE1	1:A:304:ILE:HB	2.41	0.56
1:A:106:GLU:O	1:A:110:GLN:HG2	2.06	0.56
1:I:307:GLY:HA2	2:J:63:PHE:CD1	2.41	0.56
1:C:131:ASN:ND2	1:I:243:GLY:HA3	2.21	0.56
4:I:801:SIA:H112	4:I:803:NAG:H83	1.88	0.56
1:A:29:ASN:OD1	3:A:602:NAG:O5	2.19	0.55
1:A:92:ASP:HB2	3:A:603:NAG:H82	1.88	0.55
1:C:48:CYS:HB3	1:C:281:CYS:O	2.06	0.55
1:K:133:ASP:HB3	1:K:158:VAL:HG23	1.88	0.55
1:A:266:ASN:H	1:A:266:ASN:ND2	2.05	0.55
2:J:132:GLU:HG2	2:J:138:PHE:CE2	2.40	0.55
1:K:244:ASP:OD2	1:K:245:LYS:N	2.37	0.55
1:G:302:HIS:CE1	1:G:304:ILE:HB	2.42	0.55
2:B:119:TYR:CE1	2:B:136:GLY:HA2	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:7:ASP:N	1:G:8:THR:CA	2.65	0.55
1:A:327:ILE:N	8:A:712:HOH:O	2.40	0.55
1:G:143:PRO:HD2	3:G:602:NAG:C6	2.36	0.55
2:F:23:GLY:HA3	2:F:36:ALA:HA	1.87	0.55
1:I:48:CYS:HB3	1:I:281:CYS:O	2.06	0.55
1:E:313:VAL:CG1	1:E:315:SER:H	2.20	0.55
2:B:117:ASN:ND2	8:B:209:HOH:O	2.02	0.55
1:G:7:ASP:N	2:H:140:PHE:CB	2.69	0.55
1:C:159:LYS:NZ	1:C:199:GLN:NE2	2.53	0.55
1:G:141:ALA:C	1:G:143:PRO:HD3	2.27	0.55
1:A:107:LEU:O	1:A:107:LEU:HD22	2.07	0.55
1:I:313:VAL:CG1	1:I:315:SER:H	2.20	0.55
1:C:133:ASP:HB3	1:C:158:VAL:HG23	1.88	0.55
1:A:44:ASN:ND2	1:A:291:ALA:H	2.04	0.55
1:C:9:LEU:HD12	2:D:26:HIS:HA	1.88	0.55
2:L:40:SER:N	8:L:207:HOH:O	2.39	0.55
2:D:159:TYR:H	2:D:160:PRO:CD	2.19	0.55
1:A:11:ILE:CA	2:B:14:TRP:HH2	2.20	0.55
1:C:46:LYS:CD	8:C:719:HOH:O	2.55	0.55
1:A:155:ILE:HD11	1:A:258:ARG:HD2	1.89	0.55
1:G:311:LYS:HG3	2:H:92:TRP:CE2	2.42	0.55
1:K:56:LEU:HA	8:K:913:HOH:O	1.86	0.55
1:C:119:ARG:HG2	1:C:119:ARG:O	2.04	0.55
2:H:132:GLU:HG2	2:H:138:PHE:CE2	2.40	0.55
2:L:131:LYS:CG	8:L:225:HOH:O	2.55	0.55
1:G:44:ASN:ND2	1:G:291:ALA:H	2.05	0.55
1:C:188:PRO:HG2	1:C:194:GLN:NE2	2.22	0.55
1:E:114:VAL:CG2	1:E:117:PHE:HB2	2.35	0.55
2:J:79:ASN:HB3	2:L:68:LYS:HE3	1.89	0.55
1:E:302:HIS:CE1	1:E:304:ILE:HB	2.41	0.55
1:A:297:PRO:HG3	2:B:56:ILE:HA	1.88	0.55
1:G:90:SER:CB	8:G:719:HOH:O	2.43	0.55
1:E:51:ARG:NH2	8:E:703:HOH:O	2.39	0.55
1:A:133:ASP:HB3	1:A:158:VAL:HG23	1.89	0.55
1:K:239:LEU:HD13	1:K:265:ARG:NH1	2.22	0.55
1:C:44:ASN:HD21	1:C:291:ALA:H	1.53	0.55
1:E:266:ASN:HB2	8:E:714:HOH:O	2.06	0.55
1:G:7:ASP:CB	2:H:140:PHE:HB2	2.38	0.54
2:D:132:GLU:HG2	2:D:138:PHE:CE2	2.39	0.54
1:K:48:CYS:HB3	1:K:281:CYS:O	2.06	0.54
1:E:155:ILE:HD11	1:E:258:ARG:HD2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:GLY:HA3	1:A:233:MET:O	2.08	0.54
1:K:63:ILE:HD11	1:K:85:ILE:HG21	1.90	0.54
1:C:14:HIS:HB2	2:D:20:GLY:O	2.07	0.54
1:I:99:GLY:HA3	1:I:233:MET:O	2.08	0.54
1:I:266:ASN:H	1:I:266:ASN:ND2	2.05	0.54
1:C:8:THR:OG1	2:D:27:GLN:HB3	2.07	0.54
1:K:114:VAL:CG2	1:K:117:PHE:HB2	2.36	0.54
2:L:155:GLY:N	8:L:227:HOH:O	2.34	0.54
1:A:324:LEU:HD23	1:A:324:LEU:N	2.21	0.54
1:E:44:ASN:ND2	1:E:291:ALA:H	2.04	0.54
1:I:155:ILE:HD11	1:I:258:ARG:HD2	1.89	0.54
1:E:99:GLY:HA3	1:E:233:MET:O	2.07	0.54
2:L:119:TYR:CE1	2:L:136:GLY:HA2	2.42	0.54
1:C:59:GLY:C	8:C:712:HOH:O	2.43	0.54
2:B:47:GLU:HB3	1:E:26:LEU:HB3	1.90	0.54
1:K:188:PRO:HG2	1:K:194:GLN:NE2	2.22	0.54
1:G:313:VAL:CG1	1:G:315:SER:H	2.19	0.54
1:C:313:VAL:CG1	1:C:315:SER:H	2.20	0.54
1:A:311:LYS:HE2	2:B:61:THR:HG22	1.87	0.54
1:A:189:SER:HB2	1:A:222:ILE:HD13	1.90	0.54
2:L:155:GLY:CA	8:L:227:HOH:O	2.55	0.54
1:K:313:VAL:CG1	1:K:315:SER:H	2.20	0.54
1:I:141:ALA:C	1:I:143:PRO:HD3	2.27	0.54
2:J:119:TYR:CE1	2:J:136:GLY:HA2	2.43	0.54
1:K:189:SER:HB2	1:K:222:ILE:HD13	1.90	0.54
1:G:189:SER:HB2	1:G:222:ILE:HD13	1.90	0.54
1:K:121:GLU:CD	8:K:906:HOH:O	2.45	0.54
1:A:62:ASN:HA	1:A:87:GLU:OE1	2.07	0.54
1:C:72:GLU:HG2	3:C:601:NAG:H82	1.89	0.54
2:B:95:ASN:HD21	2:D:95:ASN:ND2	2.05	0.54
1:K:141:ALA:C	1:K:143:PRO:HD3	2.28	0.54
1:A:183:TRP:HZ3	1:A:238:THR:HG22	1.73	0.54
1:I:116:SER:OG	1:I:264:GLU:CB	2.56	0.54
1:I:96:CYS:O	1:I:227:ARG:HD3	2.08	0.54
1:G:327:ILE:HD12	2:H:12:GLY:HA2	1.89	0.54
1:E:39:LEU:HB2	1:E:318:LEU:HB2	1.90	0.54
1:I:29:ASN:ND2	8:I:901:HOH:O	2.22	0.54
2:B:132:GLU:HG2	2:B:138:PHE:CE2	2.40	0.54
1:K:122:ILE:CB	8:K:914:HOH:O	2.52	0.54
1:K:324:LEU:HD23	1:K:324:LEU:N	2.23	0.54
1:K:25:VAL:HG22	2:L:102:LEU:HD12	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:SER:HA	8:A:711:HOH:O	2.06	0.54
1:E:35:SER:OG	1:E:319:ARG:NH1	2.41	0.54
1:C:222:ILE:HG23	1:C:230:GLU:CG	2.38	0.54
2:H:119:TYR:CE1	2:H:136:GLY:HA2	2.43	0.54
1:C:141:ALA:C	1:C:143:PRO:HD3	2.28	0.54
1:G:48:CYS:HB3	1:G:281:CYS:O	2.08	0.54
1:E:133:ASP:HB3	1:E:158:VAL:HG23	1.89	0.54
1:G:13:TYR:CD2	2:H:6:ILE:HG12	2.43	0.54
1:C:316:THR:HG23	1:C:317:LYS:N	2.22	0.54
1:G:312:TYR:HD2	2:H:89:LEU:HD13	1.73	0.54
1:K:14:HIS:NE2	1:K:15:ALA:C	2.61	0.53
2:H:66:VAL:C	8:H:205:HOH:O	2.11	0.53
1:I:44:ASN:HD21	1:I:291:ALA:H	1.56	0.53
1:C:172:ILE:HG22	8:C:722:HOH:O	2.06	0.53
1:C:63:ILE:HD11	1:C:85:ILE:HG21	1.90	0.53
1:G:183:TRP:HZ3	1:G:238:THR:HG22	1.73	0.53
1:A:48:CYS:HB3	1:A:281:CYS:O	2.08	0.53
1:G:155:ILE:HD11	1:G:258:ARG:HD2	1.91	0.53
1:C:314:LYS:NZ	2:F:60:ASN:O	2.42	0.53
1:G:266:ASN:H	1:G:266:ASN:ND2	2.05	0.53
1:C:44:ASN:ND2	1:C:291:ALA:H	2.06	0.53
1:E:141:ALA:C	1:E:143:PRO:HD3	2.28	0.53
1:E:96:CYS:O	1:E:227:ARG:HD3	2.08	0.53
1:E:324:LEU:N	1:E:324:LEU:HD23	2.23	0.53
1:I:190:THR:O	1:I:193:ASP:N	2.42	0.53
1:C:64:ALA:CB	8:C:729:HOH:O	2.54	0.53
2:B:127:LYS:H	2:B:127:LYS:CD	2.19	0.53
1:C:99:GLY:HA3	1:C:233:MET:O	2.09	0.53
1:C:183:TRP:HZ3	1:C:238:THR:HG22	1.73	0.53
1:K:99:GLY:HA3	1:K:233:MET:O	2.08	0.53
1:K:292:ILE:CD1	1:K:301:ILE:HD12	2.39	0.53
1:A:141:ALA:C	1:A:143:PRO:HD3	2.29	0.53
1:A:312:TYR:CD2	2:B:89:LEU:HD13	2.43	0.53
1:E:297:PRO:HG3	2:F:56:ILE:HA	1.90	0.53
1:I:44:ASN:ND2	1:I:291:ALA:H	2.07	0.53
1:E:266:ASN:C	8:E:714:HOH:O	2.46	0.53
1:A:133:ASP:CG	1:A:136:LYS:HD3	2.28	0.53
1:A:214:LYS:HG2	1:A:216:PHE:CE2	2.43	0.53
2:F:119:TYR:CE1	2:F:136:GLY:HA2	2.42	0.53
1:I:63:ILE:HD11	1:I:85:ILE:HG21	1.90	0.53
2:J:79:ASN:CB	2:L:68:LYS:HE3	2.39	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:ASN:CA	1:A:87:GLU:OE1	2.57	0.53
1:K:75:SER:N	8:K:905:HOH:O	2.36	0.53
1:K:96:CYS:O	1:K:227:ARG:HD3	2.09	0.53
1:G:99:GLY:HA3	1:G:233:MET:O	2.09	0.53
1:G:107:LEU:HD22	1:G:111:LEU:HD22	1.91	0.53
1:I:183:TRP:HZ3	1:I:238:THR:HG22	1.74	0.53
1:G:63:ILE:HD11	1:G:85:ILE:HG21	1.90	0.53
1:A:96:CYS:O	1:A:227:ARG:HD3	2.09	0.53
1:I:33:THR:HG22	1:I:326:ASN:OD1	2.09	0.53
2:D:9:PHE:CD1	2:D:10:ILE:HG13	2.44	0.53
1:K:14:HIS:O	1:K:15:ALA:HB2	2.09	0.52
2:B:76:ARG:NH1	2:D:69:GLU:O	2.40	0.52
2:F:159:TYR:CD1	2:F:161:LYS:HE3	2.43	0.52
1:K:259:TYR:C	8:K:914:HOH:O	2.44	0.52
1:I:305:THR:H	2:J:66:VAL:CG1	2.12	0.52
1:E:63:ILE:HD11	1:E:85:ILE:HG21	1.91	0.52
1:C:154:LEU:HA	8:C:717:HOH:O	2.08	0.52
2:F:9:PHE:CD1	2:F:10:ILE:HG13	2.44	0.52
1:C:96:CYS:O	1:C:227:ARG:HD3	2.09	0.52
1:I:324:LEU:N	1:I:324:LEU:HD23	2.24	0.52
1:C:302:HIS:CE1	1:C:304:ILE:HB	2.43	0.52
1:K:183:TRP:HZ3	1:K:238:THR:HG22	1.74	0.52
1:E:189:SER:HB2	1:E:222:ILE:HD13	1.91	0.52
1:G:324:LEU:HD23	1:G:324:LEU:N	2.24	0.52
2:F:132:GLU:HG2	2:F:138:PHE:CE2	2.39	0.52
1:I:312:TYR:HD2	2:J:89:LEU:HD13	1.75	0.52
1:E:281:CYS:CB	8:E:709:HOH:O	2.41	0.52
1:G:315:SER:HB3	8:H:207:HOH:O	2.08	0.52
1:E:72:GLU:CG	6:E:601:NAG:H82	2.34	0.52
1:G:113:SER:HB2	1:G:269:SER:HB2	1.92	0.52
1:C:266:ASN:ND2	1:C:266:ASN:H	2.06	0.52
1:I:50:LEU:HD13	2:J:63:PHE:CZ	2.45	0.52
1:C:311:LYS:HG3	2:D:92:TRP:CE2	2.45	0.52
1:E:312:TYR:HD2	2:F:89:LEU:HD13	1.74	0.52
1:A:62:ASN:HB2	8:A:709:HOH:O	2.09	0.52
1:I:327:ILE:N	1:I:327:ILE:HD13	2.25	0.52
1:A:159:LYS:HZ1	1:A:199:GLN:NE2	2.08	0.52
2:B:13:GLY:HA2	8:B:202:HOH:O	2.09	0.52
1:E:317:LYS:O	1:E:318:LEU:HD23	2.10	0.52
1:C:177:LYS:HB3	1:C:177:LYS:HZ2	1.75	0.52
1:C:189:SER:HB2	1:C:222:ILE:HD13	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:305:THR:HG22	2:L:66:VAL:HG12	1.91	0.52
1:I:40:GLU:OE1	1:I:294:THR:CB	2.58	0.52
1:G:7:ASP:N	2:H:140:PHE:HD1	2.08	0.51
1:A:62:ASN:N	1:A:62:ASN:OD1	2.30	0.51
2:J:106:ARG:HG3	8:J:206:HOH:O	2.10	0.51
1:G:312:TYR:CD2	2:H:89:LEU:HD13	2.45	0.51
1:C:77:SER:O	1:C:79:ALA:N	2.44	0.51
1:A:63:ILE:HD11	1:A:85:ILE:HG21	1.91	0.51
2:L:18:VAL:N	8:L:226:HOH:O	2.42	0.51
2:J:120:GLU:OE1	2:J:123:ARG:NH1	2.44	0.51
1:A:102:ILE:HG12	1:A:236:TYR:CE2	2.45	0.51
1:G:162:ASN:HD22	1:G:199:GLN:HE21	1.57	0.51
1:C:307:GLY:HA2	2:D:63:PHE:CE1	2.45	0.51
1:K:50:LEU:HD21	1:K:306:ILE:HG22	1.92	0.51
1:G:25:VAL:HG12	2:J:51:LYS:HG3	1.92	0.51
1:C:155:ILE:HG22	1:C:157:LEU:HD13	1.93	0.51
1:K:82:TRP:CB	8:K:913:HOH:O	2.55	0.51
1:G:55:PRO:HB3	1:G:84:TYR:CE2	2.46	0.51
2:H:143:LYS:HE2	2:H:143:LYS:HA	1.93	0.51
8:H:213:HOH:O	2:J:67:GLY:N	2.04	0.51
1:I:120:PHE:HA	8:I:924:HOH:O	2.06	0.51
1:E:195:GLN:HA	1:E:198:TYR:O	2.11	0.51
1:C:159:LYS:CD	1:C:199:GLN:HG2	2.40	0.51
2:H:113:SER:CB	2:L:2:LEU:HD22	2.40	0.51
1:E:183:TRP:HZ3	1:E:238:THR:HG22	1.74	0.51
1:A:223:ARG:HG3	1:C:206:PHE:HZ	1.76	0.51
2:L:9:PHE:CD1	2:L:10:ILE:HG13	2.46	0.51
1:C:44:ASN:HD22	1:C:44:ASN:C	2.13	0.51
1:E:63:ILE:HG12	1:E:87:GLU:OE1	2.11	0.51
2:H:9:PHE:CD1	2:H:10:ILE:HG13	2.46	0.51
1:I:224:PRO:O	1:I:232:ARG:NH2	2.43	0.51
1:E:312:TYR:CD2	2:F:89:LEU:HD13	2.46	0.51
1:A:55:PRO:HB3	1:A:84:TYR:CE2	2.46	0.51
2:L:29:GLU:HB2	8:L:212:HOH:O	2.10	0.50
1:G:307:GLY:HA2	2:H:63:PHE:CE1	2.46	0.50
1:E:13:TYR:CZ	2:F:6:ILE:HG23	2.46	0.50
1:I:231:GLY:C	1:I:232:ARG:HG2	2.32	0.50
1:E:55:PRO:HB3	1:E:84:TYR:CE2	2.46	0.50
2:J:25:HIS:HB2	2:J:34:TYR:CD2	2.46	0.50
1:K:162:ASN:HA	1:K:199:GLN:HE21	1.77	0.50
1:E:33:THR:HG23	1:E:324:LEU:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:ASN:C	1:A:132:HIS:CG	2.84	0.50
2:D:139:GLU:CG	8:D:212:HOH:O	2.59	0.50
2:B:28:ASN:HD22	2:B:28:ASN:H	1.60	0.50
2:L:139:GLU:OE2	8:L:211:HOH:O	2.20	0.50
1:C:63:ILE:HG12	1:C:87:GLU:OE1	2.12	0.50
1:A:159:LYS:HD3	1:A:199:GLN:H	1.77	0.50
1:G:93:ASN:ND2	3:G:602:NAG:O7	2.44	0.50
1:K:27:GLU:HB2	8:K:921:HOH:O	2.10	0.50
1:A:58:LEU:HD21	1:A:63:ILE:HD13	1.94	0.50
1:I:223:ARG:HG3	1:K:206:PHE:HZ	1.77	0.50
1:G:7:ASP:N	2:H:140:PHE:CD1	2.80	0.50
1:K:25:VAL:HG21	2:L:102:LEU:HD12	1.93	0.50
2:D:122:VAL:O	2:D:125:GLN:HB2	2.11	0.50
2:B:9:PHE:CD1	2:B:10:ILE:HG13	2.46	0.50
1:C:167:LEU:HD12	1:C:167:LEU:C	2.32	0.50
1:K:239:LEU:CD1	1:K:265:ARG:HH11	2.24	0.50
1:I:311:LYS:HE2	2:J:61:THR:HG22	1.94	0.50
1:G:122:ILE:HG23	1:G:123:PHE:N	2.27	0.50
1:I:55:PRO:HB3	1:I:84:TYR:CE2	2.47	0.50
2:B:87:GLY:HA3	2:D:88:PHE:CZ	2.47	0.50
2:B:14:TRP:CE3	2:B:17:MET:HE2	2.36	0.50
2:J:151:SER:HB2	2:J:156:THR:O	2.12	0.50
2:L:143:LYS:HE2	2:L:143:LYS:HA	1.94	0.50
1:C:74:GLU:OE1	1:C:144:HIS:CE1	2.64	0.49
1:A:159:LYS:HZ2	1:A:199:GLN:HE21	1.57	0.49
1:A:33:THR:OG1	1:A:34:HIS:ND1	2.30	0.49
1:E:311:LYS:HG3	2:F:92:TRP:CE2	2.47	0.49
1:I:190:THR:C	1:I:192:ALA:N	2.66	0.49
1:K:15:ALA:HB1	1:K:327:ILE:O	2.12	0.49
1:I:40:GLU:OE1	1:I:294:THR:HB	2.12	0.49
1:C:206:PHE:HD2	1:C:215:LYS:HB2	1.77	0.49
2:F:143:LYS:HA	2:F:143:LYS:HE2	1.93	0.49
1:G:167:LEU:C	1:G:167:LEU:HD12	2.33	0.49
1:A:33:THR:HB	1:A:34:HIS:CE1	2.47	0.49
2:J:9:PHE:CD1	2:J:10:ILE:HG13	2.47	0.49
2:J:28:ASN:H	2:J:28:ASN:HD22	1.60	0.49
1:E:282:ASN:C	1:E:282:ASN:ND2	2.66	0.49
1:E:162:ASN:HD22	1:E:199:GLN:HE21	1.59	0.49
1:K:44:ASN:C	1:K:44:ASN:HD22	2.15	0.49
4:I:801:SIA:H6	4:I:801:SIA:O1A	2.11	0.49
1:E:266:ASN:H	1:E:266:ASN:ND2	2.09	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:63:ILE:HG12	1:G:87:GLU:OE1	2.13	0.49
1:G:40:GLU:OE1	8:G:718:HOH:O	2.20	0.49
2:B:143:LYS:HA	2:B:143:LYS:HE2	1.93	0.49
2:D:143:LYS:HA	2:D:143:LYS:HE2	1.93	0.49
2:L:28:ASN:H	2:L:28:ASN:HD22	1.60	0.49
2:L:151:SER:HB2	2:L:156:THR:O	2.13	0.49
2:B:27:GLN:HG3	2:B:27:GLN:O	2.13	0.49
1:I:116:SER:OG	1:I:264:GLU:HB2	2.13	0.49
1:A:305:THR:HG22	1:A:306:ILE:N	2.27	0.49
1:K:55:PRO:HB3	1:K:84:TYR:CE2	2.47	0.49
2:D:38:LEU:O	2:D:42:GLN:HB2	2.12	0.49
1:A:12:GLY:HA3	2:B:14:TRP:CZ2	2.47	0.49
2:L:27:GLN:HG3	2:L:27:GLN:O	2.12	0.49
1:C:25:VAL:HG21	2:D:102:LEU:HD12	1.94	0.49
2:D:9:PHE:CE1	2:D:10:ILE:HG13	2.48	0.49
2:H:126:LEU:HB3	2:H:157:TYR:CE2	2.48	0.49
2:H:151:SER:HB2	2:H:156:THR:O	2.12	0.49
1:A:305:THR:CG2	1:A:306:ILE:N	2.76	0.49
1:I:116:SER:OG	1:I:264:GLU:HB3	2.13	0.49
5:E:603:SIA:H6	5:E:603:SIA:O1A	2.13	0.49
1:G:155:ILE:HG13	1:G:258:ARG:HB2	1.94	0.49
1:C:55:PRO:HB3	1:C:84:TYR:CE2	2.47	0.49
2:J:126:LEU:HB3	2:J:157:TYR:CE2	2.47	0.49
1:A:122:ILE:HG23	1:A:123:PHE:N	2.28	0.49
1:K:122:ILE:CG2	8:K:914:HOH:O	2.18	0.48
1:G:33:THR:HG23	1:G:324:LEU:O	2.13	0.48
1:A:155:ILE:HG13	1:A:258:ARG:HB2	1.95	0.48
1:E:155:ILE:HG13	1:E:258:ARG:HB2	1.95	0.48
2:F:151:SER:HB2	2:F:156:THR:O	2.13	0.48
2:J:132:GLU:CD	8:J:212:HOH:O	2.45	0.48
1:C:92:ASP:HB2	3:C:601:NAG:O6	2.12	0.48
1:I:14:HIS:NE2	8:I:909:HOH:O	2.34	0.48
1:K:305:THR:HG22	2:L:66:VAL:CG1	2.43	0.48
1:G:44:ASN:C	1:G:44:ASN:HD22	2.15	0.48
1:A:44:ASN:C	1:A:44:ASN:HD22	2.16	0.48
1:E:122:ILE:HG23	1:E:123:PHE:N	2.28	0.48
2:H:28:ASN:HD22	2:H:28:ASN:H	1.61	0.48
1:I:167:LEU:HD12	1:I:167:LEU:C	2.33	0.48
2:B:123:ARG:CG	2:B:123:ARG:O	2.62	0.48
1:A:62:ASN:CB	1:A:87:GLU:OE1	2.61	0.48
2:L:126:LEU:HB3	2:L:157:TYR:CE2	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:44:ASN:O	1:C:46:LYS:HG3	2.13	0.48
1:I:308:LYS:HE2	2:J:61:THR:O	2.12	0.48
1:K:63:ILE:HG12	1:K:87:GLU:OE1	2.12	0.48
2:B:151:SER:HB2	2:B:156:THR:O	2.12	0.48
2:L:124:SER:N	8:L:205:HOH:O	2.45	0.48
1:A:55:PRO:HB3	1:A:84:TYR:CZ	2.49	0.48
1:K:47:LEU:HD22	1:K:276:THR:O	2.14	0.48
2:H:158:ASP:C	2:H:160:PRO:HD3	2.34	0.48
1:I:76:LEU:O	1:I:77:SER:CB	2.52	0.48
1:C:176:GLY:C	1:C:177:LYS:CG	2.82	0.48
2:H:159:TYR:N	2:H:160:PRO:HD3	2.28	0.48
2:B:126:LEU:HB3	2:B:157:TYR:CE2	2.49	0.48
2:J:159:TYR:H	2:J:160:PRO:CD	2.25	0.48
2:J:162:TYR:O	8:J:216:HOH:O	2.20	0.48
1:A:159:LYS:NZ	1:A:199:GLN:HE21	2.12	0.48
2:F:9:PHE:CE1	2:F:10:ILE:HG13	2.49	0.48
1:G:113:SER:HB2	1:G:269:SER:CB	2.43	0.48
1:I:55:PRO:HB3	1:I:84:TYR:CZ	2.49	0.48
2:B:125:GLN:NE2	2:B:157:TYR:HB3	2.28	0.48
2:F:28:ASN:HD22	2:F:28:ASN:H	1.60	0.48
2:F:126:LEU:HB3	2:F:157:TYR:CE2	2.48	0.48
2:B:93:THR:O	2:B:97:GLU:HG2	2.14	0.48
2:B:124:SER:O	2:B:127:LYS:HE3	2.13	0.48
1:G:58:LEU:HD21	1:G:63:ILE:HD13	1.95	0.48
1:I:33:THR:HG23	1:I:324:LEU:O	2.14	0.48
1:G:25:VAL:HG21	2:H:102:LEU:HD22	1.95	0.48
1:K:121:GLU:OE2	8:K:906:HOH:O	2.20	0.48
2:L:124:SER:O	2:L:127:LYS:HE3	2.13	0.48
2:J:97:GLU:HB3	2:L:58:LYS:HD2	1.95	0.48
2:D:28:ASN:H	2:D:28:ASN:HD22	1.61	0.48
2:J:143:LYS:HE2	2:J:143:LYS:HA	1.94	0.48
1:A:74:GLU:CA	8:A:737:HOH:O	2.58	0.48
1:A:137:GLY:HA3	1:A:156:TRP:HB3	1.95	0.48
1:A:188:PRO:HG2	1:A:194:GLN:HE21	1.78	0.48
1:E:55:PRO:HB3	1:E:84:TYR:CZ	2.49	0.48
1:C:55:PRO:HB3	1:C:84:TYR:CZ	2.49	0.48
1:I:128:SER:N	8:I:911:HOH:O	2.47	0.48
2:D:151:SER:HB2	2:D:156:THR:O	2.13	0.48
1:A:109:GLU:N	8:A:734:HOH:O	2.45	0.47
2:J:142:HIS:NE2	2:J:161:LYS:HD3	2.24	0.47
1:E:11:ILE:HD13	2:F:119:TYR:HA	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:THR:OG1	1:A:323:GLY:HA3	2.14	0.47
1:A:188:PRO:HD3	1:A:253:ASN:ND2	2.29	0.47
2:D:126:LEU:HB3	2:D:157:TYR:CE2	2.49	0.47
2:L:17:MET:HG2	8:L:226:HOH:O	2.13	0.47
1:K:122:ILE:HG23	1:K:123:PHE:N	2.29	0.47
1:E:25:VAL:HG21	2:F:102:LEU:HD12	1.94	0.47
1:I:155:ILE:HG13	1:I:258:ARG:HB2	1.96	0.47
1:A:312:TYR:HD2	2:B:89:LEU:HD13	1.78	0.47
1:C:122:ILE:HG23	1:C:123:PHE:N	2.28	0.47
1:C:216:PHE:N	1:C:216:PHE:CD1	2.82	0.47
2:H:124:SER:O	2:H:127:LYS:HE3	2.14	0.47
2:F:124:SER:O	2:F:127:LYS:HE3	2.13	0.47
1:E:9:LEU:N	2:F:138:PHE:O	2.42	0.47
1:G:55:PRO:HB3	1:G:84:TYR:CZ	2.49	0.47
1:C:167:LEU:HD12	1:C:167:LEU:O	2.14	0.47
1:I:167:LEU:HD12	1:I:167:LEU:O	2.14	0.47
2:H:87:GLY:HA3	2:J:88:PHE:CZ	2.49	0.47
2:H:27:GLN:HG3	2:H:27:GLN:O	2.15	0.47
2:J:124:SER:O	2:J:127:LYS:HE3	2.15	0.47
1:C:223:ARG:O	1:C:230:GLU:CG	2.62	0.47
1:A:313:VAL:HG12	1:A:315:SER:N	2.28	0.47
1:A:34:HIS:N	1:A:34:HIS:ND1	2.61	0.47
1:C:129:TRP:CG	1:C:157:LEU:HD21	2.49	0.47
1:E:44:ASN:C	1:E:44:ASN:HD22	2.16	0.47
1:I:122:ILE:HG23	1:I:123:PHE:N	2.29	0.47
1:K:13:TYR:O	2:L:14:TRP:N	2.44	0.47
1:K:155:ILE:HG13	1:K:258:ARG:HB2	1.97	0.47
1:C:217:LYS:N	8:C:727:HOH:O	2.06	0.47
2:L:17:MET:SD	2:L:23:GLY:HA3	2.55	0.47
1:A:12:GLY:N	2:B:14:TRP:CZ3	2.82	0.47
2:D:161:LYS:O	2:D:162:TYR:HB2	2.15	0.47
1:A:141:ALA:O	1:A:227:ARG:NH1	2.42	0.47
1:C:117:PHE:CD2	1:C:263:MET:HG3	2.49	0.47
1:A:33:THR:HG1	1:A:34:HIS:CE1	2.28	0.47
2:B:9:PHE:CE1	2:B:10:ILE:HG13	2.50	0.47
1:E:14:HIS:CD2	1:E:15:ALA:N	2.82	0.47
2:L:9:PHE:CE1	2:L:10:ILE:HG13	2.49	0.47
8:K:925:HOH:O	2:L:13:GLY:HA2	2.15	0.47
1:K:55:PRO:HB3	1:K:84:TYR:CZ	2.50	0.47
1:A:195:GLN:HG3	1:A:196:SER:N	2.30	0.47
1:K:44:ASN:O	1:K:46:LYS:HG3	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:ASN:O	1:A:132:HIS:CD2	2.68	0.47
1:A:107:LEU:HA	1:A:107:LEU:HD23	1.69	0.47
1:C:226:VAL:HG12	1:C:227:ARG:HG3	1.97	0.47
1:G:44:ASN:O	1:G:46:LYS:HG3	2.15	0.47
1:A:173:ASN:ND2	1:A:179:VAL:HG23	2.30	0.47
1:K:13:TYR:O	2:L:13:GLY:CA	2.62	0.47
1:A:12:GLY:CA	2:B:14:TRP:CH2	2.98	0.47
1:I:44:ASN:HD21	1:I:291:ALA:N	2.12	0.47
2:D:159:TYR:O	2:D:161:LYS:N	2.48	0.47
1:A:195:GLN:CG	8:A:704:HOH:O	2.54	0.47
2:F:27:GLN:O	2:F:27:GLN:HG3	2.15	0.47
1:I:226:VAL:HG12	1:I:227:ARG:HG3	1.97	0.47
1:C:14:HIS:CD2	1:C:15:ALA:N	2.83	0.47
1:C:311:LYS:HE2	2:D:61:THR:HG22	1.96	0.47
1:K:14:HIS:CD2	1:K:15:ALA:O	2.65	0.46
1:A:132:HIS:HE1	1:A:165:PRO:O	1.98	0.46
1:K:188:PRO:HG2	1:K:194:GLN:HE21	1.80	0.46
1:G:316:THR:N	8:H:207:HOH:O	2.46	0.46
1:G:188:PRO:HG2	1:G:194:GLN:HE21	1.80	0.46
1:E:125:LYS:HB2	1:E:258:ARG:NH1	2.30	0.46
1:G:14:HIS:CD2	1:G:15:ALA:N	2.83	0.46
1:K:39:LEU:HB2	1:K:318:LEU:HB2	1.96	0.46
1:K:132:HIS:HE1	1:K:165:PRO:O	1.98	0.46
1:E:132:HIS:HE1	1:E:165:PRO:O	1.98	0.46
1:C:141:ALA:O	1:C:227:ARG:NH1	2.41	0.46
1:A:311:LYS:HG3	2:B:92:TRP:CE2	2.50	0.46
2:D:87:GLY:HA3	2:F:88:PHE:CZ	2.50	0.46
1:K:162:ASN:OD1	1:K:199:GLN:CD	2.51	0.46
1:G:125:LYS:HB2	1:G:258:ARG:NH1	2.29	0.46
1:C:177:LYS:HZ3	1:C:177:LYS:HG2	1.43	0.46
2:J:27:GLN:HG3	2:J:27:GLN:O	2.15	0.46
1:K:213:SER:N	8:K:910:HOH:O	2.38	0.46
2:H:9:PHE:CE1	2:H:10:ILE:HG13	2.50	0.46
1:G:268:GLY:N	8:G:709:HOH:O	2.48	0.46
2:B:28:ASN:H	2:B:28:ASN:ND2	2.14	0.46
1:A:11:ILE:C	2:B:14:TRP:HH2	2.18	0.46
1:A:198:TYR:O	1:A:199:GLN:C	2.54	0.46
2:D:27:GLN:HG3	2:D:27:GLN:O	2.14	0.46
1:E:25:VAL:HG22	2:F:102:LEU:HD12	1.98	0.46
1:A:44:ASN:O	1:A:46:LYS:HG3	2.16	0.46
1:G:197:LEU:HD11	7:G:603:SIA:H91	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:132:HIS:HE1	1:I:165:PRO:O	1.99	0.46
2:D:71:ASN:OD1	2:D:73:LEU:N	2.39	0.46
1:I:171:TYR:CE2	1:I:173:ASN:HA	2.51	0.46
1:A:164:TYR:CZ	1:A:252:GLY:HA2	2.51	0.46
1:K:33:THR:HG23	1:K:324:LEU:O	2.14	0.46
1:C:51:ARG:HB3	1:C:51:ARG:NH1	2.30	0.46
1:C:74:GLU:OE1	1:C:144:HIS:HE1	1.99	0.46
1:I:125:LYS:HB2	1:I:258:ARG:NH1	2.31	0.46
1:I:26:LEU:HD23	2:L:47:GLU:OE1	2.16	0.46
1:E:159:LYS:HD2	1:E:199:GLN:HG2	1.98	0.45
1:K:14:HIS:CD2	1:K:15:ALA:CA	2.98	0.45
1:E:244:ASP:OD2	1:E:245:LYS:N	2.48	0.45
1:A:22:VAL:HG21	1:A:321:ALA:HB2	1.98	0.45
1:G:7:ASP:HB2	2:H:140:PHE:HB2	1.97	0.45
2:L:28:ASN:H	2:L:28:ASN:ND2	2.14	0.45
2:B:91:ILE:HD13	2:D:91:ILE:HG21	1.98	0.45
1:E:8:THR:HG22	2:F:139:GLU:CA	2.46	0.45
1:K:226:VAL:HG12	1:K:227:ARG:HG3	1.98	0.45
1:E:188:PRO:HG2	1:E:194:GLN:HE21	1.79	0.45
1:G:167:LEU:HD12	1:G:167:LEU:O	2.15	0.45
2:J:9:PHE:CE1	2:J:10:ILE:HG13	2.51	0.45
1:E:289:LYS:HE3	1:E:289:LYS:HB3	1.40	0.45
2:H:120:GLU:OE1	2:H:123:ARG:NH1	2.49	0.45
1:K:10:CYS:O	2:L:24:TYR:HB3	2.17	0.45
1:E:226:VAL:HG12	1:E:227:ARG:HG3	1.97	0.45
2:H:91:ILE:HD13	2:J:91:ILE:HG21	1.97	0.45
1:C:132:HIS:HE1	1:C:165:PRO:O	2.00	0.45
1:I:44:ASN:HD21	1:I:291:ALA:CA	2.29	0.45
1:G:106:GLU:OE1	2:L:75:LYS:N	2.50	0.45
1:C:316:THR:HG23	1:C:317:LYS:HG2	1.97	0.45
1:G:122:ILE:HG23	1:G:123:PHE:H	1.81	0.45
1:K:311:LYS:HE2	2:L:61:THR:HG22	1.98	0.45
2:J:17:MET:SD	2:J:23:GLY:HA3	2.56	0.45
1:E:266:ASN:N	1:E:266:ASN:ND2	2.64	0.45
2:D:28:ASN:H	2:D:28:ASN:ND2	2.15	0.45
1:I:10:CYS:HA	2:J:137:CYS:HA	1.99	0.45
1:K:129:TRP:CD2	1:K:157:LEU:HD11	2.52	0.45
1:K:254:LEU:N	8:K:917:HOH:O	2.48	0.45
1:A:39:LEU:HB2	1:A:318:LEU:HB2	1.99	0.45
1:K:266:ASN:HD22	1:K:266:ASN:N	1.94	0.45
1:A:74:GLU:N	8:A:737:HOH:O	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:ASN:N	1:A:266:ASN:ND2	2.64	0.45
1:I:145:ALA:CA	8:I:903:HOH:O	2.54	0.45
1:C:159:LYS:HZ2	1:C:199:GLN:NE2	1.96	0.45
1:K:7:ASP:HA	2:L:28:ASN:HA	1.97	0.45
2:F:28:ASN:H	2:F:28:ASN:ND2	2.15	0.45
2:D:53:ASN:O	2:D:57:GLU:HB2	2.17	0.45
1:K:195:GLN:HG3	1:K:196:SER:N	2.31	0.45
1:G:164:TYR:CZ	1:G:252:GLY:HA2	2.52	0.45
1:C:39:LEU:HB2	1:C:318:LEU:HB2	1.98	0.45
1:I:164:TYR:CZ	1:I:252:GLY:HA2	2.52	0.45
1:K:106:GLU:CB	8:K:915:HOH:O	2.04	0.45
2:B:51:LYS:HA	1:E:25:VAL:O	2.16	0.45
2:H:28:ASN:ND2	2:H:28:ASN:H	2.15	0.45
2:D:87:GLY:HA3	2:F:88:PHE:CE1	2.52	0.45
1:E:327:ILE:HD12	2:F:12:GLY:HA2	1.98	0.45
1:I:28:LYS:HG3	2:L:50:ASN:OD1	2.16	0.45
1:C:214:LYS:HZ2	1:C:214:LYS:HG2	1.67	0.45
1:E:164:TYR:CZ	1:E:252:GLY:HA2	2.52	0.45
2:J:18:VAL:C	8:J:211:HOH:O	2.43	0.45
1:K:141:ALA:O	1:K:227:ARG:NH1	2.41	0.45
2:H:17:MET:SD	2:H:23:GLY:HA3	2.57	0.45
1:K:22:VAL:HG21	1:K:321:ALA:HB2	1.98	0.45
2:D:147:THR:HA	2:D:150:GLU:HB2	1.99	0.44
1:G:266:ASN:N	1:G:266:ASN:ND2	2.65	0.44
2:J:28:ASN:H	2:J:28:ASN:ND2	2.15	0.44
1:K:22:VAL:HG12	1:K:319:ARG:HG2	1.98	0.44
1:G:39:LEU:HB2	1:G:318:LEU:HB2	1.98	0.44
1:K:13:TYR:CE2	2:L:10:ILE:HG22	2.52	0.44
1:K:324:LEU:CD2	2:L:21:TRP:CD1	2.98	0.44
2:F:17:MET:SD	2:F:23:GLY:HA3	2.57	0.44
2:F:28:ASN:HD22	2:F:28:ASN:N	2.16	0.44
1:G:29:ASN:OD1	3:G:601:NAG:N2	2.50	0.44
1:G:78:THR:O	1:G:79:ALA:C	2.55	0.44
1:C:22:VAL:HG21	1:C:321:ALA:HB2	1.99	0.44
1:A:14:HIS:CD2	1:A:15:ALA:N	2.85	0.44
1:G:114:VAL:CG1	1:G:117:PHE:HB2	2.40	0.44
1:E:54:ALA:HB1	1:E:55:PRO:HD2	1.98	0.44
1:K:164:TYR:CZ	1:K:252:GLY:HA2	2.52	0.44
1:I:195:GLN:HG3	1:I:196:SER:N	2.31	0.44
1:A:76:LEU:CA	8:A:738:HOH:O	2.48	0.44
1:A:33:THR:CB	1:A:34:HIS:ND1	2.79	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:159:LYS:HD2	1:G:199:GLN:HG2	1.98	0.44
2:B:23:GLY:HA3	2:B:36:ALA:HA	2.00	0.44
1:I:51:ARG:HB2	1:I:53:VAL:CG1	2.48	0.44
1:K:54:ALA:HB1	1:K:55:PRO:HD2	2.00	0.44
1:K:266:ASN:O	1:K:267:ALA:C	2.55	0.44
1:A:226:VAL:HG12	1:A:227:ARG:HG3	1.99	0.44
1:C:231:GLY:N	8:C:701:HOH:O	2.50	0.44
2:B:87:GLY:HA3	2:D:88:PHE:CE1	2.53	0.44
1:K:47:LEU:HD12	1:K:286:GLN:NE2	2.32	0.44
1:E:22:VAL:HG21	1:E:321:ALA:HB2	2.00	0.44
1:A:7:ASP:HA	1:A:8:THR:HA	1.79	0.44
1:A:108:ARG:CB	8:A:734:HOH:O	1.88	0.44
1:K:125:LYS:HB2	1:K:258:ARG:NH1	2.33	0.44
1:A:72:GLU:C	1:A:74:GLU:H	2.21	0.44
2:H:160:PRO:HD2	2:H:161:LYS:HG3	2.00	0.44
1:C:164:TYR:CZ	1:C:252:GLY:HA2	2.52	0.44
1:C:305:THR:HG22	2:D:66:VAL:HG12	1.99	0.44
1:K:51:ARG:CZ	1:K:51:ARG:HB3	2.48	0.44
1:K:105:GLU:OE2	8:K:920:HOH:O	2.21	0.44
1:C:14:HIS:HB2	2:D:21:TRP:HA	2.00	0.44
1:I:53:VAL:HG22	1:I:53:VAL:O	2.18	0.44
1:E:129:TRP:CD2	1:E:157:LEU:HD11	2.53	0.44
1:G:22:VAL:HG21	1:G:321:ALA:HB2	1.99	0.44
1:C:188:PRO:HG2	1:C:194:GLN:HE21	1.81	0.44
1:K:311:LYS:HG3	2:L:92:TRP:CE2	2.52	0.44
2:J:125:GLN:HE22	2:J:155:GLY:C	2.21	0.44
1:E:307:GLY:HA2	2:F:63:PHE:CE1	2.52	0.44
2:J:123:ARG:HB2	2:J:138:PHE:HZ	1.83	0.43
2:B:14:TRP:CE3	2:B:17:MET:HE1	2.53	0.43
2:L:131:LYS:HE3	2:L:133:ILE:HD13	2.00	0.43
2:J:28:ASN:N	2:J:28:ASN:HD22	2.16	0.43
1:E:205:VAL:HG21	1:E:254:LEU:HD13	2.00	0.43
2:H:30:GLN:H	2:H:30:GLN:HE21	1.66	0.43
1:K:122:ILE:HG23	1:K:123:PHE:H	1.82	0.43
1:C:76:LEU:HB3	1:C:77:SER:H	1.47	0.43
1:A:233:MET:HE3	1:A:233:MET:HB2	1.91	0.43
1:A:204:TYR:CD1	1:A:204:TYR:O	2.71	0.43
1:I:265:ARG:NH2	8:I:910:HOH:O	2.50	0.43
2:B:160:PRO:HD2	2:B:161:LYS:HG3	2.00	0.43
1:A:122:ILE:HG23	1:A:123:PHE:H	1.83	0.43
1:G:145:ALA:C	1:G:147:ALA:H	2.22	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:605:SIA:H6	4:A:605:SIA:O1A	2.17	0.43
1:I:72:GLU:C	1:I:74:GLU:H	2.22	0.43
1:E:313:VAL:HG12	1:E:315:SER:N	2.30	0.43
1:C:22:VAL:HG12	1:C:319:ARG:HG2	1.99	0.43
2:L:159:TYR:CD1	2:L:161:LYS:HD2	2.53	0.43
1:G:72:GLU:C	1:G:74:GLU:H	2.21	0.43
1:K:13:TYR:CZ	2:L:12:GLY:O	2.71	0.43
1:A:302:HIS:HA	1:A:303:PRO:HD3	1.84	0.43
1:I:314:LYS:NZ	2:L:60:ASN:O	2.46	0.43
1:I:41:ASP:OD1	1:I:41:ASP:C	2.56	0.43
2:B:71:ASN:ND2	2:B:71:ASN:N	2.30	0.43
1:I:313:VAL:HG12	1:I:315:SER:N	2.30	0.43
2:D:28:ASN:HD22	2:D:28:ASN:N	2.16	0.43
2:B:30:GLN:HE21	2:B:30:GLN:H	1.66	0.43
2:F:125:GLN:HE22	2:F:155:GLY:C	2.21	0.43
1:K:14:HIS:O	2:L:13:GLY:HA3	2.18	0.43
1:K:16:ASN:N	8:K:904:HOH:O	2.30	0.43
1:A:159:LYS:HZ1	1:A:199:GLN:HE22	1.67	0.43
2:D:120:GLU:O	2:D:124:SER:OG	2.37	0.43
1:C:195:GLN:HG3	1:C:196:SER:N	2.31	0.43
1:A:214:LYS:HE3	1:A:216:PHE:CE2	2.53	0.43
1:C:107:LEU:O	1:C:108:ARG:C	2.55	0.43
1:G:54:ALA:HB1	1:G:55:PRO:HD2	2.00	0.43
1:E:195:GLN:HG3	1:E:196:SER:N	2.31	0.43
1:E:122:ILE:HG23	1:E:123:PHE:H	1.83	0.43
1:C:173:ASN:ND2	1:C:179:VAL:HG23	2.33	0.43
1:I:12:GLY:N	2:J:14:TRP:CH2	2.87	0.43
1:I:42:LYS:HE3	1:I:42:LYS:HB2	1.87	0.43
2:J:159:TYR:O	2:J:161:LYS:N	2.52	0.43
2:L:30:GLN:HE21	2:L:30:GLN:H	1.67	0.43
1:C:117:PHE:HD2	1:C:263:MET:HG3	1.82	0.43
1:E:319:ARG:NH1	1:E:319:ARG:HG3	2.30	0.43
1:E:47:LEU:HD22	1:E:276:THR:O	2.19	0.43
1:E:296:LEU:HD23	1:E:296:LEU:HA	1.70	0.43
1:G:311:LYS:HE2	2:H:61:THR:HG22	2.00	0.43
1:K:276:THR:HA	1:K:277:PRO:HD3	1.87	0.43
1:C:122:ILE:HG23	1:C:123:PHE:H	1.83	0.43
1:E:173:ASN:ND2	1:E:179:VAL:HG23	2.34	0.43
1:I:39:LEU:HB2	1:I:318:LEU:HB2	2.01	0.43
2:F:30:GLN:H	2:F:30:GLN:HE21	1.67	0.43
2:J:26:HIS:NE2	2:J:31:GLY:O	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:313:VAL:HG12	1:G:315:SER:N	2.30	0.43
1:I:76:LEU:N	1:I:76:LEU:HD23	2.34	0.43
1:G:195:GLN:HG3	1:G:196:SER:N	2.32	0.43
1:I:78:THR:O	1:I:79:ALA:C	2.57	0.43
2:D:79:ASN:ND2	1:E:109:GLU:OE2	2.52	0.43
2:L:152:VAL:HG23	8:L:223:HOH:O	2.18	0.43
2:B:26:HIS:C	2:B:26:HIS:CD2	2.91	0.43
1:I:197:LEU:HD11	4:I:801:SIA:H91	2.01	0.43
1:K:307:GLY:HA2	2:L:63:PHE:CE1	2.54	0.43
2:J:30:GLN:HE21	2:J:30:GLN:H	1.67	0.43
1:C:72:GLU:C	1:C:74:GLU:H	2.21	0.42
1:A:33:THR:C	1:A:34:HIS:ND1	2.73	0.42
2:D:30:GLN:H	2:D:30:GLN:HE21	1.67	0.42
1:K:173:ASN:ND2	1:K:179:VAL:HG23	2.33	0.42
2:B:17:MET:HE2	2:B:17:MET:HB2	1.76	0.42
1:K:114:VAL:CA	1:K:266:ASN:HD21	2.32	0.42
1:G:11:ILE:O	2:H:10:ILE:HD13	2.19	0.42
1:I:51:ARG:HB3	1:I:51:ARG:NH1	2.34	0.42
2:F:71:ASN:O	8:F:203:HOH:O	2.21	0.42
1:C:273:ILE:N	1:C:273:ILE:HD12	2.33	0.42
1:G:223:ARG:HG3	1:I:206:PHE:HZ	1.84	0.42
1:C:78:THR:O	1:C:79:ALA:C	2.58	0.42
1:C:159:LYS:HD2	1:C:199:GLN:HG2	2.01	0.42
1:G:268:GLY:CA	8:G:709:HOH:O	2.59	0.42
1:C:266:ASN:ND2	1:C:266:ASN:N	2.65	0.42
1:I:22:VAL:HG21	1:I:321:ALA:HB2	2.00	0.42
1:C:11:ILE:HD11	2:D:24:TYR:HE2	1.73	0.42
1:I:145:ALA:C	1:I:147:ALA:H	2.22	0.42
1:K:72:GLU:C	1:K:74:GLU:H	2.22	0.42
1:E:318:LEU:HA	1:E:318:LEU:HD23	1.75	0.42
1:A:169:LYS:HG3	1:A:170:SER:H	1.84	0.42
2:L:5:ALA:HB3	2:L:112:ASP:OD2	2.20	0.42
1:K:167:LEU:C	1:K:167:LEU:HD23	2.40	0.42
1:A:167:LEU:HD12	1:A:168:SER:N	2.32	0.42
1:A:266:ASN:HD22	1:A:266:ASN:N	2.18	0.42
1:A:54:ALA:HB1	1:A:55:PRO:HD2	2.01	0.42
1:A:273:ILE:N	1:A:273:ILE:HD12	2.35	0.42
1:A:62:ASN:HB2	1:A:87:GLU:OE1	2.19	0.42
1:K:13:TYR:CE2	2:L:6:ILE:HA	2.53	0.42
1:E:8:THR:HA	2:F:138:PHE:O	2.20	0.42
1:A:324:LEU:H	1:A:324:LEU:HD23	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:176:GLY:C	1:I:177:LYS:HG2	2.39	0.42
1:E:58:LEU:O	1:E:76:LEU:HD11	2.19	0.42
1:C:145:ALA:C	1:C:147:ALA:H	2.23	0.42
2:F:4:GLY:O	2:F:8:GLY:HA3	2.20	0.42
1:K:145:ALA:C	1:K:147:ALA:H	2.23	0.42
2:H:75:LYS:HG3	1:I:109:GLU:HG2	2.01	0.42
1:E:305:THR:HG21	8:E:719:HOH:O	2.18	0.42
1:E:8:THR:CA	2:F:138:PHE:O	2.68	0.42
2:H:131:LYS:HE3	2:H:133:ILE:HD13	2.02	0.42
2:D:116:LYS:O	2:D:119:TYR:HB3	2.20	0.42
1:E:68:LEU:O	1:E:151:TYR:HB3	2.19	0.42
1:E:72:GLU:C	1:E:74:GLU:H	2.22	0.42
1:I:125:LYS:CB	1:I:258:ARG:NH1	2.83	0.42
1:I:229:GLN:HB3	8:I:922:HOH:O	2.19	0.42
1:K:68:LEU:O	1:K:151:TYR:HB3	2.19	0.42
1:E:273:ILE:HD12	1:E:273:ILE:N	2.34	0.42
1:A:91:SER:OG	8:A:717:HOH:O	2.21	0.42
2:J:102:LEU:O	2:J:106:ARG:HG3	2.20	0.42
1:I:244:ASP:OD2	1:I:245:LYS:N	2.53	0.42
2:L:125:GLN:HE22	2:L:155:GLY:C	2.23	0.42
1:E:8:THR:HA	2:F:139:GLU:HA	2.02	0.42
1:C:313:VAL:HG12	1:C:315:SER:N	2.30	0.42
1:G:68:LEU:O	1:G:151:TYR:HB3	2.20	0.42
2:F:127:LYS:H	2:F:127:LYS:CD	2.21	0.42
2:L:18:VAL:C	8:L:226:HOH:O	2.49	0.41
1:C:77:SER:O	1:C:78:THR:C	2.59	0.41
1:A:236:TYR:HE1	8:A:721:HOH:O	2.03	0.41
1:C:69:GLY:O	1:C:152:LYS:CG	2.62	0.41
1:K:239:LEU:HD13	1:K:265:ARG:HH11	1.84	0.41
2:F:127:LYS:N	2:F:127:LYS:HD2	2.24	0.41
2:B:106:ARG:HH21	2:F:106:ARG:NH1	2.17	0.41
1:I:50:LEU:HD13	2:J:63:PHE:HZ	1.83	0.41
1:A:125:LYS:HB2	1:A:258:ARG:NH1	2.35	0.41
1:E:233:MET:HE3	1:E:233:MET:HB2	1.91	0.41
1:G:125:LYS:CB	1:G:258:ARG:NH1	2.83	0.41
1:E:145:ALA:C	1:E:147:ALA:H	2.23	0.41
1:A:327:ILE:CB	8:A:712:HOH:O	2.68	0.41
1:A:141:ALA:O	1:A:143:PRO:HD3	2.20	0.41
1:K:313:VAL:HG12	1:K:315:SER:N	2.30	0.41
1:I:231:GLY:O	1:I:232:ARG:HG2	2.20	0.41
2:J:5:ALA:HB3	2:J:112:ASP:OD2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:173:ASN:ND2	1:G:179:VAL:HG23	2.35	0.41
2:D:131:LYS:HE3	2:D:133:ILE:HD13	2.01	0.41
2:B:59:MET:HG3	2:B:59:MET:O	2.19	0.41
1:I:190:THR:CG2	1:I:193:ASP:CA	2.98	0.41
2:B:123:ARG:HB2	2:B:138:PHE:HZ	1.84	0.41
1:A:100:ASP:O	1:A:102:ILE:CD1	2.68	0.41
1:C:70:ASN:HA	1:C:71:PRO:HD3	1.96	0.41
2:B:4:GLY:O	2:B:8:GLY:HA3	2.19	0.41
1:I:122:ILE:HG23	1:I:123:PHE:H	1.84	0.41
1:G:144:HIS:HE1	8:G:716:HOH:O	2.03	0.41
2:F:131:LYS:HE3	2:F:133:ILE:HD13	2.02	0.41
1:C:51:ARG:HB3	1:C:51:ARG:HH11	1.86	0.41
1:C:244:ASP:OD2	1:C:245:LYS:N	2.54	0.41
2:B:91:ILE:HG21	2:F:91:ILE:HD13	2.01	0.41
1:E:13:TYR:HE2	2:F:6:ILE:HA	1.78	0.41
1:I:125:LYS:HB2	1:I:258:ARG:HH11	1.85	0.41
1:G:269:SER:N	8:G:702:HOH:O	1.85	0.41
1:A:244:ASP:OD2	1:A:245:LYS:N	2.54	0.41
1:G:273:ILE:N	1:G:273:ILE:HD12	2.36	0.41
1:K:161:GLY:O	1:K:162:ASN:HB3	2.20	0.41
1:C:222:ILE:CG2	1:C:230:GLU:HG2	2.46	0.41
1:I:13:TYR:HB2	1:I:324:LEU:HD11	2.03	0.41
1:K:144:HIS:O	1:K:145:ALA:HB3	2.21	0.41
2:B:131:LYS:HE3	2:B:133:ILE:HD13	2.02	0.41
1:E:167:LEU:C	1:E:167:LEU:HD23	2.41	0.41
1:G:51:ARG:NH1	1:G:51:ARG:HB3	2.35	0.41
1:C:258:ARG:HH11	1:C:258:ARG:HD2	1.72	0.41
1:E:51:ARG:NH1	1:E:51:ARG:HB3	2.35	0.41
1:K:266:ASN:O	1:K:267:ALA:O	2.38	0.41
1:I:200:ASN:OD1	1:I:200:ASN:N	2.54	0.41
1:K:302:HIS:HA	1:K:303:PRO:HD3	1.85	0.41
1:G:25:VAL:CG1	2:J:51:LYS:HG3	2.51	0.41
1:I:70:ASN:HA	1:I:71:PRO:HD3	1.93	0.41
1:A:8:THR:CG2	2:B:138:PHE:C	2.80	0.41
1:I:141:ALA:O	1:I:227:ARG:NH1	2.42	0.41
1:A:160:LYS:O	1:A:161:GLY:C	2.59	0.41
1:G:125:LYS:HB2	1:G:258:ARG:HH11	1.84	0.41
1:E:311:LYS:HE2	2:F:61:THR:HG22	2.02	0.41
1:E:118:GLU:CG	1:E:262:ALA:HB3	2.51	0.41
1:G:118:GLU:CG	1:G:262:ALA:HB3	2.50	0.41
1:K:297:PRO:HG3	2:L:56:ILE:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:159:TYR:N	2:D:160:PRO:HD2	2.36	0.41
1:A:114:VAL:O	8:A:740:HOH:O	2.22	0.41
1:A:144:HIS:O	1:A:145:ALA:HB3	2.21	0.41
1:I:25:VAL:HG11	2:L:51:LYS:HG3	1.98	0.41
1:C:13:TYR:HA	2:D:21:TRP:O	2.20	0.41
2:H:28:ASN:HD22	2:H:28:ASN:N	2.16	0.41
2:J:87:GLY:HA3	2:L:88:PHE:CZ	2.56	0.41
1:A:327:ILE:CA	8:A:712:HOH:O	2.69	0.41
1:C:77:SER:C	1:C:79:ALA:N	2.74	0.41
1:K:292:ILE:CG2	1:K:294:THR:HG22	2.51	0.41
1:E:294:THR:HG22	8:E:704:HOH:O	2.20	0.41
1:G:296:LEU:HA	1:G:297:PRO:HD3	1.87	0.41
1:G:162:ASN:HA	1:G:162:ASN:HD22	1.74	0.41
1:E:141:ALA:O	1:E:227:ARG:NH1	2.41	0.41
1:E:125:LYS:CB	1:E:258:ARG:NH1	2.83	0.41
1:I:26:LEU:HG	2:L:47:GLU:HB3	2.03	0.41
1:I:41:ASP:OD1	1:I:41:ASP:N	2.49	0.41
2:J:131:LYS:HE3	2:J:133:ILE:HD13	2.03	0.41
2:H:4:GLY:O	2:H:8:GLY:HA3	2.21	0.41
1:A:70:ASN:HA	1:A:71:PRO:HD3	1.95	0.41
2:H:127:LYS:HD2	2:H:127:LYS:N	2.23	0.41
2:B:4:GLY:HA2	8:B:212:HOH:O	2.12	0.41
2:H:51:LYS:HG3	1:K:25:VAL:HG12	2.02	0.40
1:E:63:ILE:O	1:E:67:ILE:HG12	2.21	0.40
1:C:302:HIS:HA	1:C:303:PRO:HD3	1.85	0.40
2:D:139:GLU:HG3	8:D:212:HOH:O	2.18	0.40
1:G:14:HIS:N	2:H:21:TRP:O	2.54	0.40
1:I:68:LEU:O	1:I:151:TYR:HB3	2.20	0.40
1:E:107:LEU:HB2	1:E:237:TRP:CE2	2.57	0.40
1:C:276:THR:HA	1:C:277:PRO:HD3	1.87	0.40
1:C:157:LEU:HD12	1:C:157:LEU:HA	1.81	0.40
1:K:125:LYS:HB2	1:K:258:ARG:HH11	1.87	0.40
1:K:71:PRO:O	1:K:74:GLU:HB2	2.21	0.40
1:I:118:GLU:CG	1:I:262:ALA:HB3	2.51	0.40
1:G:13:TYR:CZ	2:H:6:ILE:HG23	2.56	0.40
1:I:63:ILE:O	1:I:67:ILE:HG12	2.22	0.40
1:G:155:ILE:HG22	1:G:157:LEU:HD13	2.03	0.40
1:G:118:GLU:HG3	1:G:262:ALA:HB3	2.03	0.40
1:A:51:ARG:HH21	1:A:53:VAL:HG21	1.86	0.40
1:A:68:LEU:O	1:A:151:TYR:HB3	2.21	0.40
2:L:4:GLY:O	2:L:8:GLY:HA3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:173:ASN:HD21	1:I:179:VAL:HG23	1.87	0.40
1:C:54:ALA:HB1	1:C:55:PRO:HD2	2.03	0.40
1:G:244:ASP:OD2	1:G:245:LYS:N	2.55	0.40
1:K:119:ARG:HB2	1:K:261:PHE:CD2	2.56	0.40
1:K:42:LYS:HB2	1:K:42:LYS:HE3	1.88	0.40
1:E:308:LYS:HE3	1:E:308:LYS:HB2	1.93	0.40
1:I:273:ILE:N	1:I:273:ILE:HD12	2.36	0.40
2:B:97:GLU:CA	8:B:213:HOH:O	2.24	0.40
2:D:161:LYS:O	2:D:162:TYR:CB	2.70	0.40
2:L:157:TYR:HD1	8:L:223:HOH:O	1.89	0.40
1:K:24:THR:HG22	2:L:104:ASN:CB	2.47	0.40
1:C:63:ILE:O	1:C:67:ILE:HG12	2.21	0.40
1:G:129:TRP:CG	1:G:157:LEU:HD21	2.57	0.40
2:J:87:GLY:HA3	2:L:88:PHE:CE1	2.57	0.40
1:I:204:TYR:CD1	1:I:215:LYS:HE3	2.57	0.40
1:E:78:THR:O	1:E:79:ALA:C	2.59	0.40
1:K:273:ILE:HD12	1:K:273:ILE:N	2.36	0.40
1:I:120:PHE:N	8:I:924:HOH:O	2.50	0.40
1:E:266:ASN:N	1:E:266:ASN:HD22	2.18	0.40
2:J:136:GLY:HA3	8:J:208:HOH:O	2.20	0.40
1:G:63:ILE:O	1:G:67:ILE:HG12	2.21	0.40
2:D:119:TYR:CE1	2:D:136:GLY:HA2	2.56	0.40
1:E:204:TYR:CD1	1:E:215:LYS:HE3	2.56	0.40

All (4) 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:E:119:ARG:NH2	2:L:38:LEU:CD2[1_565]	1.96	0.24
1:E:79:ALA:N	2:L:150:GLU:OE2[1_565]	2.01	0.19
1:E:258:ARG:NH2	2:L:42:GLN:OE1[1_565]	2.04	0.16
1:E:81:SER:OG	2:L:146:ASN:OD1[1_565]	2.12	0.08

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	319/321 (99%)	295 (92%)	23 (7%)	1 (0%)	46	84
1	C	319/321 (99%)	296 (93%)	19 (6%)	4 (1%)	15	53
1	E	319/321 (99%)	295 (92%)	23 (7%)	1 (0%)	46	84
1	G	319/321 (99%)	292 (92%)	27 (8%)	0	100	100
1	I	319/321 (99%)	291 (91%)	26 (8%)	2 (1%)	30	72
1	K	319/321 (99%)	290 (91%)	29 (9%)	0	100	100
2	B	160/162 (99%)	146 (91%)	13 (8%)	1 (1%)	30	72
2	D	160/162 (99%)	143 (89%)	16 (10%)	1 (1%)	30	72
2	F	159/162 (98%)	145 (91%)	14 (9%)	0	100	100
2	H	160/162 (99%)	146 (91%)	14 (9%)	0	100	100
2	J	160/162 (99%)	144 (90%)	16 (10%)	0	100	100
2	L	159/162 (98%)	145 (91%)	13 (8%)	1 (1%)	30	72
All	All	2872/2898 (99%)	2628 (92%)	233 (8%)	11 (0%)	39	80

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	160	PRO
1	A	146	GLY
1	C	77	SER
1	C	78	THR
1	C	161	GLY
1	I	78	THR
1	C	79	ALA
2	L	161	LYS
2	D	119	TYR
1	E	79	ALA
1	I	79	ALA

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	281/282 (100%)	248 (88%)	33 (12%)	7	27
1	C	282/282 (100%)	248 (88%)	34 (12%)	6	25
1	E	282/282 (100%)	254 (90%)	28 (10%)	10	35
1	G	281/282 (100%)	258 (92%)	23 (8%)	14	46
1	I	282/282 (100%)	247 (88%)	35 (12%)	6	24
1	K	282/282 (100%)	255 (90%)	27 (10%)	10	38
2	B	139/139 (100%)	128 (92%)	11 (8%)	15	48
2	D	138/139 (99%)	125 (91%)	13 (9%)	11	39
2	F	139/139 (100%)	126 (91%)	13 (9%)	11	39
2	H	139/139 (100%)	130 (94%)	9 (6%)	21	58
2	J	138/139 (99%)	128 (93%)	10 (7%)	18	53
2	L	139/139 (100%)	128 (92%)	11 (8%)	15	48
All	All	2522/2526 (100%)	2275 (90%)	247 (10%)	10	36

All (247) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	THR
1	A	13	TYR
1	A	27	GLU
1	A	34	HIS
1	A	37	ASN
1	A	44	ASN
1	A	50	LEU
1	A	58	LEU
1	A	90	SER
1	A	101	PHE
1	A	102	ILE
1	A	107	LEU
1	A	111	LEU
1	A	136	LYS
1	A	157	LEU
1	A	160	LYS
1	A	167	LEU
1	A	169	LYS
1	A	177	LYS
1	A	190	THR

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Mol	Chain	Res	Type
1	A	195	GLN
1	A	196	SER
1	A	198	TYR
1	A	211	ARG
1	A	228	ASP
1	A	240	VAL
1	A	265	ARG
1	A	266	ASN
1	A	275	ASP
1	A	316	THR
1	A	324	LEU
1	A	326	ASN
1	A	327	ILE
2	B	24	TYR
2	B	28	ASN
2	B	30	GLN
2	B	38	LEU
2	B	68	LYS
2	B	71	ASN
2	B	80	LEU
2	B	106	ARG
2	B	123	ARG
2	B	126	LEU
2	B	154	ASN
1	C	26	LEU
1	C	37	ASN
1	C	44	ASN
1	C	51	ARG
1	C	53	VAL
1	C	67	ILE
1	C	90	SER
1	C	107	LEU
1	C	111	LEU
1	C	115	SER
1	C	119	ARG
1	C	148	LYS
1	C	152	LYS
1	C	160	LYS
1	C	167	LEU
1	C	168	SER
1	C	177	LYS
1	C	190	THR

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Mol	Chain	Res	Type
1	C	195	GLN
1	C	198	TYR
1	C	211	ARG
1	C	214	LYS
1	C	216	PHE
1	C	228	ASP
1	C	230	GLU
1	C	258	ARG
1	C	263	MET
1	C	265	ARG
1	C	266	ASN
1	C	305	THR
1	C	316	THR
1	C	317	LYS
1	C	324	LEU
1	C	325	ARG
2	D	17	MET
2	D	22	TYR
2	D	28	ASN
2	D	30	GLN
2	D	38	LEU
2	D	66	VAL
2	D	80	LEU
2	D	98	LEU
2	D	124	SER
2	D	126	LEU
2	D	150	GLU
2	D	154	ASN
2	D	158	ASP
1	E	9	LEU
1	E	32	VAL
1	E	37	ASN
1	E	44	ASN
1	E	47	LEU
1	E	51	ARG
1	E	67	ILE
1	E	101	PHE
1	E	111	LEU
1	E	114	VAL
1	E	167	LEU
1	E	190	THR
1	E	195	GLN

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Mol	Chain	Res	Type
1	E	198	TYR
1	E	205	VAL
1	E	211	ARG
1	E	214	LYS
1	E	217	LYS
1	E	228	ASP
1	E	230	GLU
1	E	232	ARG
1	E	266	ASN
1	E	282	ASN
1	E	289	LYS
1	E	305	THR
1	E	316	THR
1	E	319	ARG
1	E	324	LEU
2	F	2	LEU
2	F	19	ASP
2	F	22	TYR
2	F	28	ASN
2	F	30	GLN
2	F	66	VAL
2	F	80	LEU
2	F	83	LYS
2	F	98	LEU
2	F	122	VAL
2	F	126	LEU
2	F	154	ASN
2	F	159	TYR
1	G	37	ASN
1	G	44	ASN
1	G	50	LEU
1	G	51	ARG
1	G	58	LEU
1	G	67	ILE
1	G	107	LEU
1	G	111	LEU
1	G	157	LEU
1	G	167	LEU
1	G	190	THR
1	G	195	GLN
1	G	198	TYR
1	G	214	LYS

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Mol	Chain	Res	Type
1	G	228	ASP
1	G	240	VAL
1	G	265	ARG
1	G	266	ASN
1	G	269	SER
1	G	305	THR
1	G	316	THR
1	G	319	ARG
1	G	324	LEU
2	H	22	TYR
2	H	28	ASN
2	H	30	GLN
2	H	38	LEU
2	H	80	LEU
2	H	123	ARG
2	H	126	LEU
2	H	154	ASN
2	H	158	ASP
1	I	37	ASN
1	I	44	ASN
1	I	50	LEU
1	I	51	ARG
1	I	53	VAL
1	I	67	ILE
1	I	75	SER
1	I	76	LEU
1	I	77	SER
1	I	101	PHE
1	I	111	LEU
1	I	116	SER
1	I	134	SER
1	I	157	LEU
1	I	167	LEU
1	I	172	ILE
1	I	177	LYS
1	I	178	GLU
1	I	191	SER
1	I	194	GLN
1	I	195	GLN
1	I	198	TYR
1	I	200	ASN
1	I	211	ARG

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Mol	Chain	Res	Type
1	I	214	LYS
1	I	228	ASP
1	I	263	MET
1	I	265	ARG
1	I	266	ASN
1	I	289	LYS
1	I	305	THR
1	I	316	THR
1	I	319	ARG
1	I	324	LEU
1	I	327	ILE
2	J	22	TYR
2	J	26	HIS
2	J	28	ASN
2	J	30	GLN
2	J	66	VAL
2	J	80	LEU
2	J	98	LEU
2	J	126	LEU
2	J	154	ASN
2	J	158	ASP
1	K	9	LEU
1	K	11	ILE
1	K	32	VAL
1	K	37	ASN
1	K	44	ASN
1	K	47	LEU
1	K	50	LEU
1	K	51	ARG
1	K	80	SER
1	K	101	PHE
1	K	111	LEU
1	K	114	VAL
1	K	167	LEU
1	K	190	THR
1	K	195	GLN
1	K	198	TYR
1	K	205	VAL
1	K	211	ARG
1	K	214	LYS
1	K	228	ASP
1	K	246	ILE

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Mol	Chain	Res	Type
1	K	265	ARG
1	K	266	ASN
1	K	293	ASN
1	K	305	THR
1	K	316	THR
1	K	324	LEU
2	L	2	LEU
2	L	22	TYR
2	L	28	ASN
2	L	30	GLN
2	L	66	VAL
2	L	80	LEU
2	L	98	LEU
2	L	122	VAL
2	L	126	LEU
2	L	154	ASN
2	L	159	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (91) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	44	ASN
1	A	132	HIS
1	A	162	ASN
1	A	194	GLN
1	A	195	GLN
1	A	199	GLN
1	A	266	ASN
2	B	26	HIS
2	B	27	GLN
2	B	28	ASN
2	B	71	ASN
2	B	125	GLN
2	B	129	ASN
2	B	146	ASN
1	C	14	HIS
1	C	37	ASN
1	C	44	ASN
1	C	131	ASN
1	C	132	HIS
1	C	144	HIS

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Mol	Chain	Res	Type
1	C	162	ASN
1	C	194	GLN
1	C	199	GLN
1	C	253	ASN
1	C	266	ASN
2	D	27	GLN
2	D	28	ASN
2	D	95	ASN
2	D	129	ASN
2	D	142	HIS
2	D	146	ASN
1	E	14	HIS
1	E	37	ASN
1	E	44	ASN
1	E	132	HIS
1	E	162	ASN
1	E	194	GLN
1	E	199	GLN
1	E	253	ASN
1	E	266	ASN
1	E	282	ASN
2	F	27	GLN
2	F	28	ASN
2	F	125	GLN
2	F	129	ASN
2	F	146	ASN
1	G	14	HIS
1	G	37	ASN
1	G	44	ASN
1	G	162	ASN
1	G	194	GLN
1	G	199	GLN
1	G	253	ASN
1	G	266	ASN
2	H	27	GLN
2	H	28	ASN
2	H	95	ASN
2	H	129	ASN
2	H	146	ASN
1	I	37	ASN
1	I	44	ASN
1	I	93	ASN

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Mol	Chain	Res	Type
1	I	132	HIS
1	I	162	ASN
1	I	194	GLN
1	I	199	GLN
1	I	253	ASN
1	I	266	ASN
2	J	26	HIS
2	J	27	GLN
2	J	28	ASN
2	J	125	GLN
2	J	129	ASN
2	J	142	HIS
2	J	146	ASN
1	K	37	ASN
1	K	44	ASN
1	K	93	ASN
1	K	132	HIS
1	K	194	GLN
1	K	199	GLN
1	K	253	ASN
1	K	266	ASN
1	K	293	ASN
2	L	27	GLN
2	L	28	ASN
2	L	125	GLN
2	L	128	ASN
2	L	129	ASN
2	L	146	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 ⓘ

19 carbohydrates are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SIA	A	605	4	16,20,21	0.42	0	18,28,31	2.29	1 (5%)
4	GAL	A	606	4	11,11,12	0.61	0	14,15,17	0.88	0
4	NAG	A	607	4	14,14,15	0.59	0	15,19,21	0.77	0
5	SIA	C	602	5	16,20,21	0.32	0	18,28,31	2.48	1 (5%)
5	GAL	C	603	5	11,11,12	0.56	0	14,15,17	0.92	1 (7%)
5	NAG	C	604	5	14,14,15	0.56	0	15,19,21	1.40	2 (13%)
5	GAL	C	605	5	11,11,12	0.68	0	14,15,17	1.03	1 (7%)
6	NAG	E	601	1,6	14,14,15	0.55	0	15,19,21	0.81	1 (6%)
6	NAG	E	602	6	14,14,15	0.43	0	15,19,21	0.86	0
5	SIA	E	603	5	16,20,21	0.30	0	18,28,31	2.17	2 (11%)
5	GAL	E	604	5	11,11,12	0.67	0	14,15,17	0.85	1 (7%)
5	NAG	E	605	5	14,14,15	0.60	0	15,19,21	0.72	0
5	GAL	E	606	5	11,11,12	0.61	0	14,15,17	0.59	0
4	SIA	I	801	4	16,20,21	0.33	0	18,28,31	2.26	1 (5%)
4	GAL	I	802	4	11,11,12	0.70	0	14,15,17	1.26	2 (14%)
4	NAG	I	803	4	14,14,15	0.48	0	15,19,21	0.70	0
4	SIA	K	801	4	16,20,21	0.30	0	18,28,31	2.27	1 (5%)
4	GAL	K	802	4	11,11,12	0.60	0	14,15,17	0.89	0
4	NAG	K	803	4	14,14,15	0.50	0	15,19,21	0.68	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SIA	A	605	4	-	0/14/34/38	0/1/1/1
4	GAL	A	606	4	-	0/2/19/22	0/1/1/1
4	NAG	A	607	4	-	0/6/23/26	0/1/1/1
5	SIA	C	602	5	-	0/14/34/38	0/1/1/1
5	GAL	C	603	5	-	0/2/19/22	0/1/1/1
5	NAG	C	604	5	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GAL	C	605	5	-	0/2/19/22	0/1/1/1
6	NAG	E	601	1,6	-	0/6/23/26	0/1/1/1
6	NAG	E	602	6	-	0/6/23/26	0/1/1/1
5	SIA	E	603	5	-	0/14/34/38	0/1/1/1
5	GAL	E	604	5	-	0/2/19/22	0/1/1/1
5	NAG	E	605	5	-	0/6/23/26	0/1/1/1
5	GAL	E	606	5	-	0/2/19/22	0/1/1/1
4	SIA	I	801	4	-	0/14/34/38	0/1/1/1
4	GAL	I	802	4	-	0/2/19/22	0/1/1/1
4	NAG	I	803	4	-	0/6/23/26	0/1/1/1
4	SIA	K	801	4	-	0/14/34/38	0/1/1/1
4	GAL	K	802	4	-	0/2/19/22	0/1/1/1
4	NAG	K	803	4	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	602	SIA	O6-C2-C3	-10.28	90.09	109.86
4	A	605	SIA	O6-C2-C3	-9.32	91.93	109.86
4	K	801	SIA	O6-C2-C3	-9.10	92.37	109.86
4	I	801	SIA	O6-C2-C3	-9.08	92.39	109.86
5	E	603	SIA	O6-C2-C3	-8.71	93.11	109.86
4	I	802	GAL	C1-O5-C5	-2.20	109.46	112.25
5	E	603	SIA	C7-C6-C5	-2.09	111.17	114.32
5	E	604	GAL	C1-O5-C5	-2.02	109.68	112.25
6	E	601	NAG	C1-O5-C5	2.03	114.83	112.25
5	C	604	NAG	C4-C3-C2	2.26	114.74	111.23
5	C	603	GAL	O5-C5-C6	2.33	112.40	107.35
5	C	605	GAL	C3-C4-C5	2.39	114.37	110.20
4	I	802	GAL	C1-C2-C3	3.05	113.15	109.54
5	C	604	NAG	C1-O5-C5	3.22	116.33	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	605	SIA	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	606	GAL	1	0
4	A	607	NAG	2	0
5	C	602	SIA	1	0
6	E	601	NAG	3	0
6	E	602	NAG	1	0
5	E	603	SIA	2	0
4	I	801	SIA	3	0
4	I	803	NAG	1	0

5.6 Ligand geometry [i](#)

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	A	601	1	14,14,15	0.50	0	15,19,21	0.79	0
3	NAG	A	602	1	14,14,15	0.54	0	15,19,21	0.58	0
3	NAG	A	603	1	14,14,15	0.52	0	15,19,21	0.60	0
3	NAG	A	604	1	14,14,15	0.54	0	15,19,21	0.59	0
3	NAG	C	601	1	14,14,15	0.44	0	15,19,21	1.13	1 (6%)
3	NAG	G	601	1	14,14,15	0.28	0	15,19,21	0.53	0
3	NAG	G	602	1	14,14,15	0.74	0	15,19,21	1.58	2 (13%)
7	SIA	G	603	-	16,20,21	0.26	0	18,28,31	1.41	4 (22%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	601	1	-	0/6/23/26	0/1/1/1
3	NAG	A	602	1	-	0/6/23/26	0/1/1/1
3	NAG	A	603	1	-	0/6/23/26	0/1/1/1
3	NAG	A	604	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	C	601	1	-	0/6/23/26	0/1/1/1
3	NAG	G	601	1	-	0/6/23/26	0/1/1/1
3	NAG	G	602	1	-	0/6/23/26	0/1/1/1
7	SIA	G	603	-	-	0/14/34/38	0/1/1/1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	G	603	SIA	C7-C6-C5	-2.69	110.25	114.32
7	G	603	SIA	C3-C4-C5	-2.15	109.08	111.47
3	G	602	NAG	C1-O5-C5	-2.07	109.62	112.25
7	G	603	SIA	O6-C6-C5	2.98	113.36	108.48
7	G	603	SIA	O6-C2-C3	3.18	115.98	109.86
3	C	601	NAG	C1-O5-C5	3.34	116.49	112.25
3	G	602	NAG	C2-N2-C7	4.38	128.67	123.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	NAG	4	0
3	A	603	NAG	2	0
3	C	601	NAG	2	0
3	G	601	NAG	1	0
3	G	602	NAG	4	0
7	G	603	SIA	1	0

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	321/321 (100%)	0.18	12 (3%) 45 19	35, 69, 111, 154	0
1	C	321/321 (100%)	0.28	22 (6%) 20 7	36, 68, 118, 223	0
1	E	321/321 (100%)	0.18	10 (3%) 52 24	41, 71, 121, 203	0
1	G	321/321 (100%)	0.33	18 (5%) 28 11	39, 76, 124, 154	0
1	I	321/321 (100%)	0.28	12 (3%) 45 19	39, 75, 117, 169	0
1	K	321/321 (100%)	0.55	25 (7%) 16 6	52, 91, 136, 187	0
2	B	162/162 (100%)	0.38	14 (8%) 13 4	35, 84, 142, 180	0
2	D	162/162 (100%)	0.56	17 (10%) 8 3	39, 86, 163, 220	0
2	F	161/162 (99%)	0.76	29 (18%) 2 1	43, 86, 160, 246	0
2	H	162/162 (100%)	0.46	17 (10%) 8 3	40, 85, 149, 181	0
2	J	162/162 (100%)	0.32	6 (3%) 45 19	47, 80, 129, 156	0
2	L	161/162 (99%)	1.87	58 (36%) 0 0	43, 101, 223, 362	0
All	All	2896/2898 (99%)	0.44	240 (8%) 14 5	35, 78, 142, 362	0

All (240) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	129	ASN	12.5
2	L	146	ASN	9.9
2	L	149	MET	9.4
2	L	130	ALA	9.2
2	L	147	THR	8.8
2	L	154	ASN	8.5
2	L	23	GLY	8.3
2	L	159	TYR	8.3
1	G	9	LEU	8.1
1	K	77	SER	7.0
2	L	138	PHE	6.9

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Mol	Chain	Res	Type	RSRZ
2	L	160	PRO	6.7
2	L	158	ASP	6.7
1	G	75	SER	6.6
2	F	139	GLU	6.4
2	L	143	LYS	6.0
2	L	156	THR	5.8
1	K	10	CYS	5.8
2	L	153	LYS	5.7
1	G	76	LEU	5.7
2	D	129	ASN	5.7
2	L	24	TYR	5.6
2	L	133	ILE	5.6
1	E	11	ILE	5.5
1	C	74	GLU	5.5
2	L	124	SER	5.5
2	H	33	GLY	5.5
1	C	7	ASP	5.5
2	F	159	TYR	5.4
2	L	150	GLU	5.4
2	L	126	LEU	5.4
1	K	9	LEU	5.3
2	F	23	GLY	5.3
1	K	147	ALA	5.2
2	F	158	ASP	5.2
2	D	31	GLY	5.1
2	L	141	TYR	5.0
2	D	143	LYS	5.0
2	F	29	GLU	5.0
2	L	31	GLY	4.9
2	L	140	PHE	4.9
2	L	122	VAL	4.8
2	F	144	CYS	4.7
2	L	32	SER	4.7
2	D	139	GLU	4.6
2	D	158	ASP	4.6
2	L	29	GLU	4.4
2	F	128	ASN	4.4
2	L	35	ALA	4.2
2	L	125	GLN	4.2
1	K	11	ILE	4.2
1	G	211	ARG	4.1
2	L	121	LYS	4.1

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Mol	Chain	Res	Type	RSRZ
2	F	138	PHE	4.1
2	F	143	LYS	4.0
2	D	141	TYR	3.9
2	L	127	LYS	3.9
1	C	12	GLY	3.9
1	E	12	GLY	3.9
2	L	145	ASP	3.8
2	F	141	TYR	3.8
1	C	293	ASN	3.7
2	H	27	GLN	3.7
1	G	162	ASN	3.7
2	F	35	ALA	3.7
1	A	50	LEU	3.6
1	I	78	THR	3.6
2	H	134	GLY	3.6
2	F	142	HIS	3.6
2	L	42	GLN	3.6
1	G	77	SER	3.6
1	K	76	LEU	3.5
2	H	156	THR	3.5
2	L	139	GLU	3.5
1	C	8	THR	3.5
2	J	157	TYR	3.4
2	L	142	HIS	3.4
2	B	140	PHE	3.4
2	B	156	THR	3.4
1	K	50	LEU	3.4
1	K	56	LEU	3.4
2	F	30	GLN	3.4
1	I	91	SER	3.3
1	G	12	GLY	3.3
2	B	143	LYS	3.3
2	F	121	LYS	3.3
2	L	131	LYS	3.3
2	H	141	TYR	3.2
2	L	155	GLY	3.2
2	L	132	GLU	3.2
2	L	20	GLY	3.2
2	L	26	HIS	3.2
1	K	248	PHE	3.1
1	G	11	ILE	3.1
2	L	128	ASN	3.1

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Mol	Chain	Res	Type	RSRZ
1	K	8	THR	3.1
2	H	144	CYS	3.1
1	A	162	ASN	3.1
1	K	326	ASN	3.1
2	H	145	ASP	3.0
2	B	138	PHE	3.0
2	B	32	SER	3.0
1	I	93	ASN	3.0
2	L	134	GLY	3.0
1	A	19	THR	3.0
2	L	22	TYR	3.0
1	E	9	LEU	3.0
2	D	132	GLU	2.9
1	A	15	ALA	2.9
1	G	17	ASN	2.9
2	H	154	ASN	2.9
1	A	8	THR	2.9
1	I	8	THR	2.9
1	C	16	ASN	2.9
2	D	146	ASN	2.9
2	F	140	PHE	2.9
2	L	15	THR	2.9
2	L	25	HIS	2.9
2	J	156	THR	2.8
2	B	31	GLY	2.8
2	L	151	SER	2.8
2	L	36	ALA	2.8
2	H	23	GLY	2.8
1	C	195	GLN	2.8
2	H	30	GLN	2.8
2	L	54	SER	2.8
2	D	33	GLY	2.8
2	L	21	TRP	2.8
2	H	139	GLU	2.7
1	C	77	SER	2.7
2	B	147	THR	2.7
2	F	31	GLY	2.7
2	L	148	CYS	2.7
2	B	158	ASP	2.7
2	J	143	LYS	2.7
1	E	76	LEU	2.7
2	F	153	LYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	I	92	ASP	2.7
2	F	145	ASP	2.7
2	H	29	GLU	2.6
2	H	26	HIS	2.6
2	D	157	TYR	2.6
2	F	131	LYS	2.6
2	F	149	MET	2.6
2	L	119	TYR	2.6
2	D	140	PHE	2.5
2	B	128	ASN	2.5
2	L	19	ASP	2.5
1	G	8	THR	2.5
2	F	24	TYR	2.5
1	E	195	GLN	2.5
1	G	195	GLN	2.5
2	B	27	GLN	2.5
1	C	60	LYS	2.5
1	I	144	HIS	2.5
1	C	42	LYS	2.5
2	B	145	ASP	2.5
1	C	291	ALA	2.5
2	L	10	ILE	2.5
1	C	294	THR	2.5
2	L	33	GLY	2.5
1	C	228	ASP	2.5
2	J	31	GLY	2.5
1	C	51	ARG	2.4
1	K	84	TYR	2.4
1	A	133	ASP	2.4
1	K	92	ASP	2.4
2	H	159	TYR	2.4
1	G	53	VAL	2.4
2	L	34	TYR	2.4
1	G	158	VAL	2.4
1	C	76	LEU	2.4
2	D	32	SER	2.4
2	D	19	ASP	2.4
2	L	18	VAL	2.4
2	L	30	GLN	2.4
1	I	226	VAL	2.4
1	K	72	GLU	2.3
1	I	268	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	10	CYS	2.3
1	K	21	THR	2.3
1	K	140	ALA	2.3
2	B	146	ASN	2.3
2	F	125	GLN	2.3
2	F	126	LEU	2.3
1	K	17	ASN	2.2
1	K	307	GLY	2.2
2	D	11	GLU	2.2
2	F	22	TYR	2.2
2	J	131	LYS	2.2
1	C	292	ILE	2.2
2	L	135	ASN	2.2
1	I	280	ASP	2.2
1	G	293	ASN	2.2
1	E	136	LYS	2.2
2	L	9	PHE	2.2
1	A	73	CYS	2.2
2	F	160	PRO	2.2
2	H	22	TYR	2.2
1	K	228	ASP	2.2
1	K	52	GLY	2.2
2	D	149	MET	2.2
1	A	163	SER	2.1
1	E	77	SER	2.1
1	G	149	SER	2.1
1	A	292	ILE	2.1
2	D	133	ILE	2.1
1	E	13	TYR	2.1
2	H	143	LYS	2.1
1	K	29	ASN	2.1
2	H	148	CYS	2.1
1	C	147	ALA	2.1
1	C	20	ASP	2.1
1	K	7	ASP	2.1
1	C	46	LYS	2.1
1	I	94	GLY	2.1
2	B	137	CYS	2.1
2	F	27	GLN	2.1
1	G	152	LYS	2.1
2	D	18	VAL	2.1
1	I	208	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	7	ASP	2.1
1	C	41	ASP	2.1
2	F	119	TYR	2.1
1	K	43	HIS	2.0
2	B	150	GLU	2.0
1	E	162	ASN	2.0
1	K	47	LEU	2.0
1	C	268	GLY	2.0
2	J	136	GLY	2.0
1	A	201	ALA	2.0
2	F	12	GLY	2.0
1	C	10	CYS	2.0
1	E	10	CYS	2.0
2	L	11	GLU	2.0
1	G	129	TRP	2.0
1	I	198	TYR	2.0
1	G	79	ALA	2.0
1	K	93	ASN	2.0
2	F	147	THR	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	SIA	E	603	20/21	0.93	0.23	0.09	42,65,87,99	0
4	SIA	K	801	20/21	0.94	0.22	-0.41	64,83,91,91	0
4	SIA	A	605	20/21	0.94	0.18	-0.54	37,53,76,80	0
5	SIA	C	602	20/21	0.94	0.19	-0.69	48,71,90,97	0
6	NAG	E	601	14/15	0.90	0.16	-0.71	74,82,98,102	0
4	SIA	I	801	20/21	0.94	0.19	-0.72	61,77,95,96	0
4	GAL	I	802	11/12	0.88	0.22	-0.90	61,95,108,114	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	NAG	E	605	14/15	0.86	0.34	-	120,126,146,147	0
6	NAG	E	602	14/15	0.76	0.36	-	88,113,132,144	0
5	GAL	C	603	11/12	0.93	0.17	-	56,84,92,100	0
5	GAL	C	605	11/12	0.60	0.25	-	104,129,151,153	0
4	GAL	A	606	11/12	0.94	0.10	-	53,93,104,109	0
5	NAG	C	604	14/15	0.83	0.25	-	75,106,127,128	0
4	NAG	I	803	14/15	0.89	0.27	-	111,129,140,142	0
5	GAL	E	606	11/12	0.77	0.26	-	110,139,162,174	0
4	NAG	K	803	14/15	0.80	0.23	-	110,146,159,172	0
4	GAL	K	802	11/12	0.79	0.22	-	96,131,149,169	0
5	GAL	E	604	11/12	0.92	0.14	-	78,95,110,122	0
4	NAG	A	607	14/15	0.85	0.23	-	65,80,89,96	0

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	NAG	G	602	14/15	0.82	0.28	2.09	100,121,147,149	0
7	SIA	G	603	20/21	0.89	0.28	0.53	65,86,119,122	0
3	NAG	A	603	14/15	0.88	0.28	0.52	70,102,144,146	0
3	NAG	C	601	14/15	0.79	0.23	0.39	70,86,97,103	0
3	NAG	A	604	14/15	0.71	0.26	-	99,145,158,161	0
3	NAG	A	602	14/15	0.70	0.31	-	146,167,182,187	0
3	NAG	G	601	14/15	0.55	0.48	-	119,147,162,173	0
3	NAG	A	601	14/15	0.68	0.34	-	142,168,177,178	0

6.5 Other polymers [i](#)

There are no such residues in this entry.