



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 05:28 PM GMT

PDB ID : 3OGM
Title : Structure of COI1-ASK1 in complex with coronatine and the JAZ1 degron
Authors : Sheard, L.B.; Tan, X.; Mao, H.; Withers, J.; Ben-Nissan, G.; Hinds, T.R.;
Hsu, F.; Sharon, M.; Browse, J.; He, S.Y.; Rizo, J.; Howe, G.A.; Zheng, N.
Deposited on : 2010-08-17
Resolution : 3.34 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

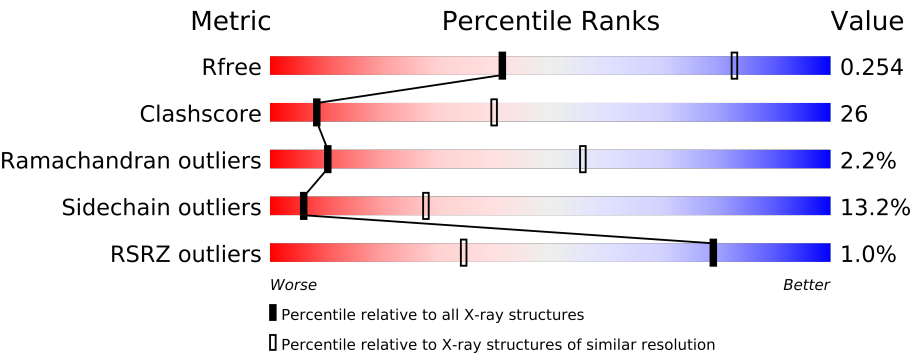
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.34 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}	66092	1421 (3.48-3.20)
Clashscore	79885	1077 (3.46-3.22)
Ramachandran outliers	78287	1053 (3.46-3.22)
Sidechain outliers	78261	1052 (3.46-3.22)
RSRZ outliers	66119	1422 (3.48-3.20)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	160	
1	C	160	
1	E	160	
1	G	160	
1	I	160	
1	K	160	
1	M	160	
1	O	160	
2	B	592	
2	D	592	
2	F	592	
2	H	592	
2	J	592	
2	L	592	

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Mol	Chain	Length	Quality of chain
2	N	592	
2	P	592	
3	Q	21	
3	R	21	
3	S	21	
3	U	21	
3	V	21	
3	W	21	
3	X	21	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
5	PO4	B	1101	-	X
5	PO4	B	1102	-	X
5	PO4	B	1103	-	X
5	PO4	B	1104	-	X
5	PO4	F	1101	-	X
5	PO4	F	1102	-	X
5	PO4	F	1103	-	X
5	PO4	H	1101	-	X
5	PO4	H	1102	-	X
5	PO4	H	1103	-	X
5	PO4	H	1104	-	X
5	PO4	J	1101	-	X
5	PO4	J	1102	-	X
5	PO4	L	1101	-	X
5	PO4	L	1102	-	X
5	PO4	L	1103	-	X
5	PO4	N	1102	-	X
5	PO4	N	1103	-	X
5	PO4	P	1101	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 46877 atoms, of which 0 are hydrogen and 0 are deuterium.

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 SKP1-like protein 1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	144	Total	C	N	O	S	0	0	0
			1146	720	185	235	6			
1	C	144	Total	C	N	O	S	0	0	0
			1146	720	185	235	6			
1	E	144	Total	C	N	O	S	0	0	0
			1146	720	185	235	6			
1	G	144	Total	C	N	O	S	0	0	0
			1146	720	185	235	6			
1	I	144	Total	C	N	O	S	0	0	0
			1146	720	185	235	6			
1	K	144	Total	C	N	O	S	0	0	0
			1146	720	185	235	6			
1	M	144	Total	C	N	O	S	0	0	0
			1146	720	185	235	6			
1	O	144	Total	C	N	O	S	0	0	0
			1146	720	185	235	6			

- Molecule 2 is a protein called Coronatine-insensitiveprotein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	568	Total	C	N	O	S	0	0	0
			4541	2873	790	842	36			
2	D	568	Total	C	N	O	S	0	0	0
			4541	2873	790	842	36			
2	F	568	Total	C	N	O	S	0	0	0
			4541	2873	790	842	36			
2	H	562	Total	C	N	O	S	0	0	0
			4486	2840	779	831	36			
2	J	568	Total	C	N	O	S	0	0	0
			4541	2873	790	842	36			
2	L	568	Total	C	N	O	S	0	0	0
			4541	2873	790	842	36			

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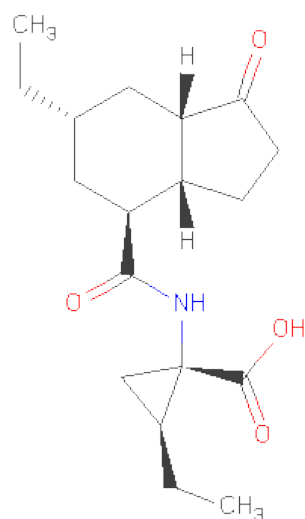
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	N	568	Total	C	N	O	S	0	0	0
			4541	2873	790	842	36			
2	P	568	Total	C	N	O	S	0	0	0
			4541	2873	790	842	36			

- Molecule 3 is a protein called JAZ1 degron peptide.

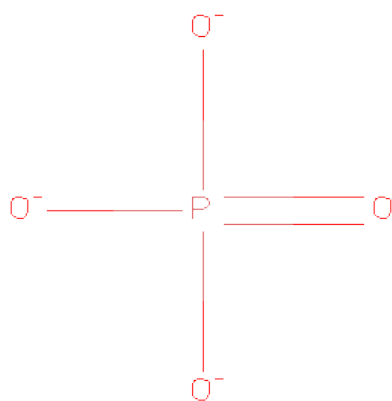
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	Q	18	Total	C	N	O	0	0	0
			156	99	34	23			
3	R	18	Total	C	N	O	0	0	0
			156	99	34	23			
3	S	18	Total	C	N	O	0	0	0
			156	99	34	23			
3	U	18	Total	C	N	O	0	0	0
			156	99	34	23			
3	V	18	Total	C	N	O	0	0	0
			156	99	34	23			
3	W	18	Total	C	N	O	0	0	0
			156	99	34	23			
3	X	18	Total	C	N	O	0	0	0
			156	99	34	23			

- Molecule 4 is (1S,2S)-2-ETHYL-1-({[(3AS,4S,6R,7AS)-6-ETHYL-1-OXOOCTAHYDRO-1H-INDEN-4-YL]CARBONYL}AMINO)CYCLOPROPANECARBOXYLICACID (three-letter code: OGK) (formula: C₁₈H₂₇NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			23	18	1	4		
4	D	1	Total	C	N	O	0	0
			23	18	1	4		
4	F	1	Total	C	N	O	0	0
			23	18	1	4		
4	H	1	Total	C	N	O	0	0
			23	18	1	4		
4	J	1	Total	C	N	O	0	0
			23	18	1	4		
4	L	1	Total	C	N	O	0	0
			23	18	1	4		
4	N	1	Total	C	N	O	0	0
			23	18	1	4		
4	P	1	Total	C	N	O	0	0
			23	18	1	4		

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	P	0	0
			5	4	1		
5	B	1	Total	O	P	0	0
			5	4	1		
5	B	1	Total	O	P	0	0
			5	4	1		
5	B	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	O	P	0	0
			5	4	1		
5	D	1	Total	O	P	0	0
			5	4	1		
5	D	1	Total	O	P	0	0
			5	4	1		
5	D	1	Total	O	P	0	0
			5	4	1		
5	F	1	Total	O	P	0	0
			5	4	1		
5	F	1	Total	O	P	0	0
			5	4	1		
5	F	1	Total	O	P	0	0
			5	4	1		
5	F	1	Total	O	P	0	0
			5	4	1		
5	H	1	Total	O	P	0	0
			5	4	1		
5	H	1	Total	O	P	0	0
			5	4	1		
5	H	1	Total	O	P	0	0
			5	4	1		
5	H	1	Total	O	P	0	0
			5	4	1		
5	J	1	Total	O	P	0	0
			5	4	1		
5	J	1	Total	O	P	0	0
			5	4	1		
5	J	1	Total	O	P	0	0
			5	4	1		
5	J	1	Total	O	P	0	0
			5	4	1		
5	L	1	Total	O	P	0	0
			5	4	1		
5	L	1	Total	O	P	0	0
			5	4	1		
5	L	1	Total	O	P	0	0
			5	4	1		
5	L	1	Total	O	P	0	0
			5	4	1		
5	N	1	Total	O	P	0	0
			5	4	1		

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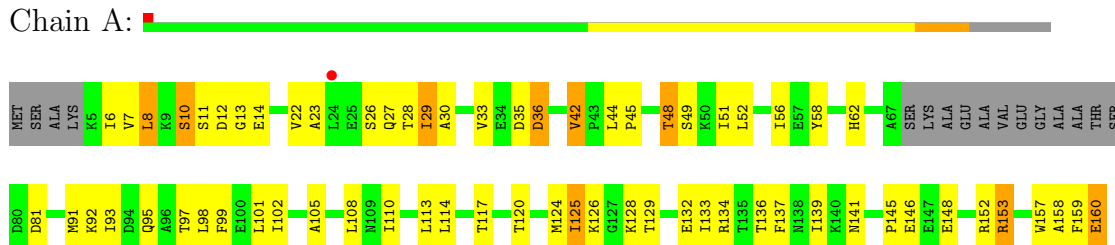
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	N	1	Total	O	P	0	0
			5	4	1		
5	N	1	Total	O	P	0	0
			5	4	1		
5	N	1	Total	O	P	0	0
			5	4	1		
5	P	1	Total	O	P	0	0
			5	4	1		
5	P	1	Total	O	P	0	0
			5	4	1		
5	P	1	Total	O	P	0	0
			5	4	1		
5	P	1	Total	O	P	0	0
			5	4	1		

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.

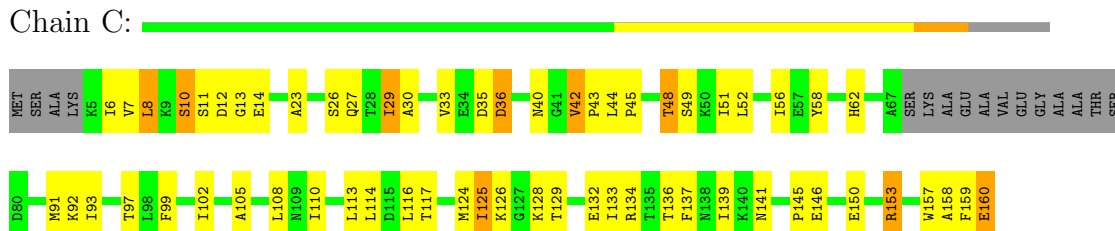
- Molecule 1: SKP1-like protein 1A

Chain A:



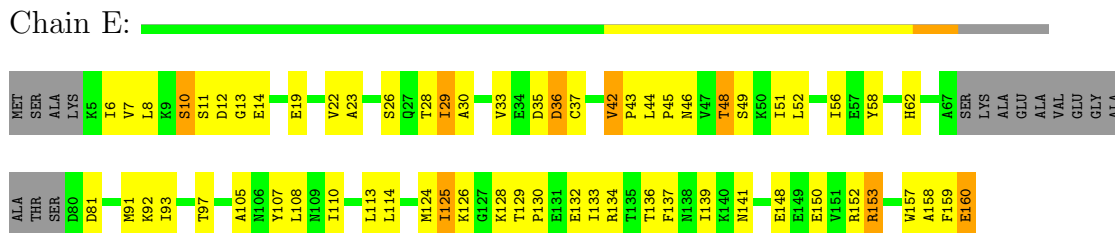
- Molecule 1: SKP1-like protein 1A

Chain C:



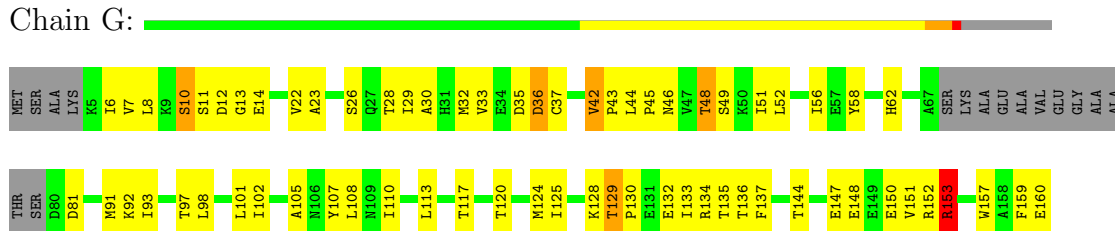
- Molecule 1: SKP1-like protein 1A

Chain E:



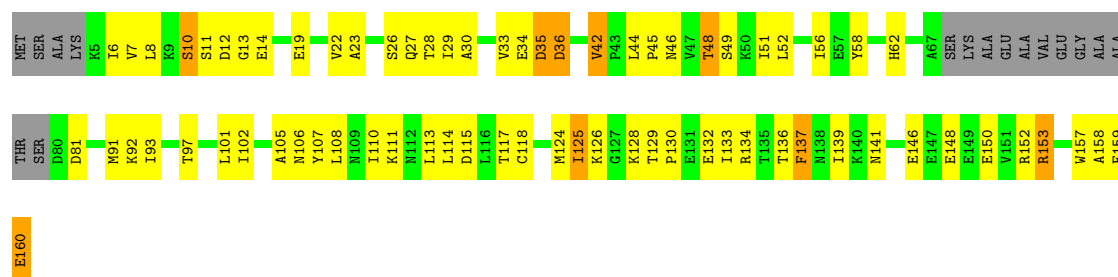
- Molecule 1: SKP1-like protein 1A

Chain G:



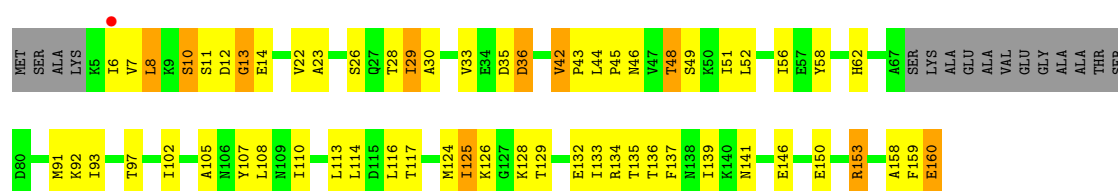
- Molecule 1: SKP1-like protein 1A

Chain I:



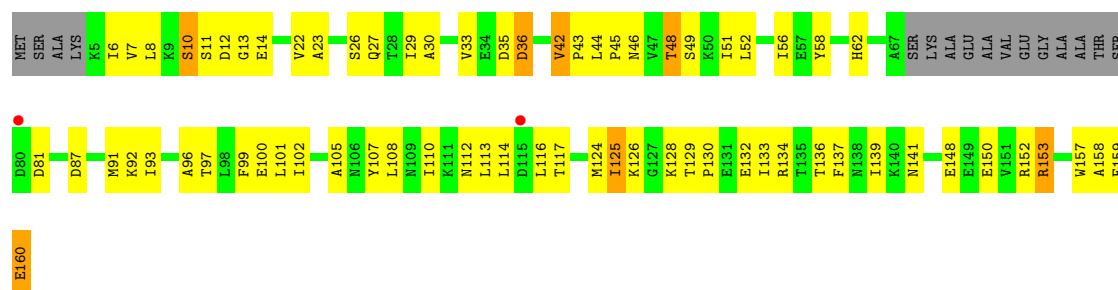
- Molecule 1: SKP1-like protein 1A

Chain K:



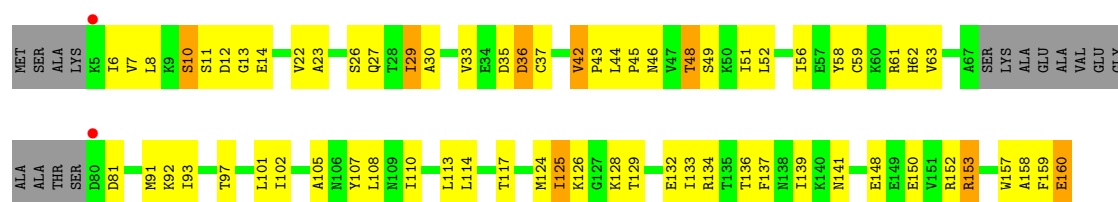
- Molecule 1: SKP1-like protein 1A

Chain M:



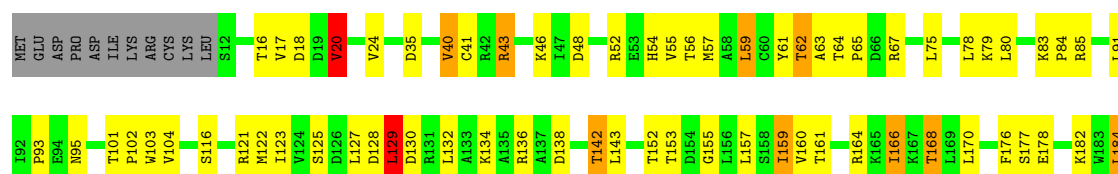
- Molecule 1: SKP1-like protein 1A

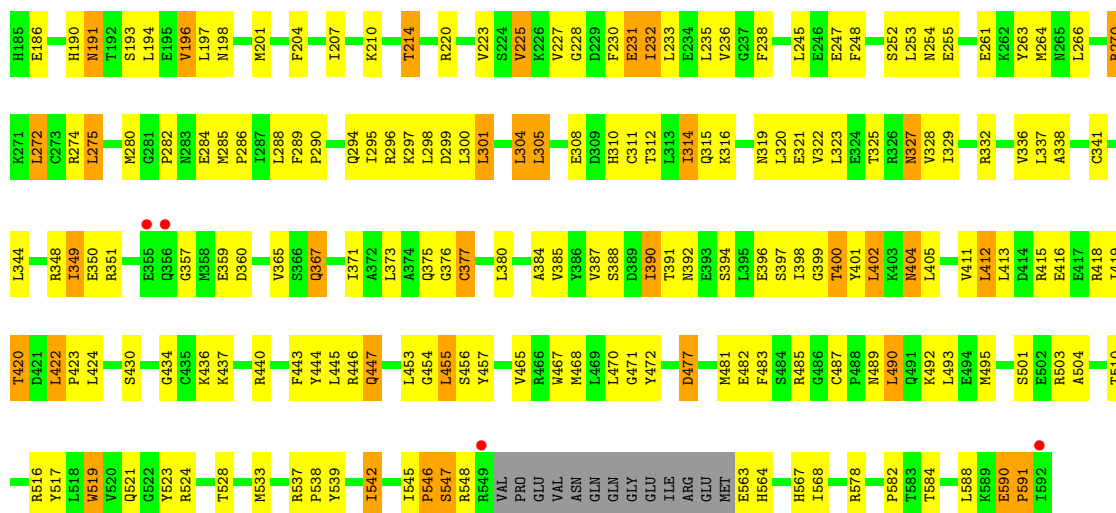
Chain O:



- Molecule 2: Coronatine-insensitiveprotein 1

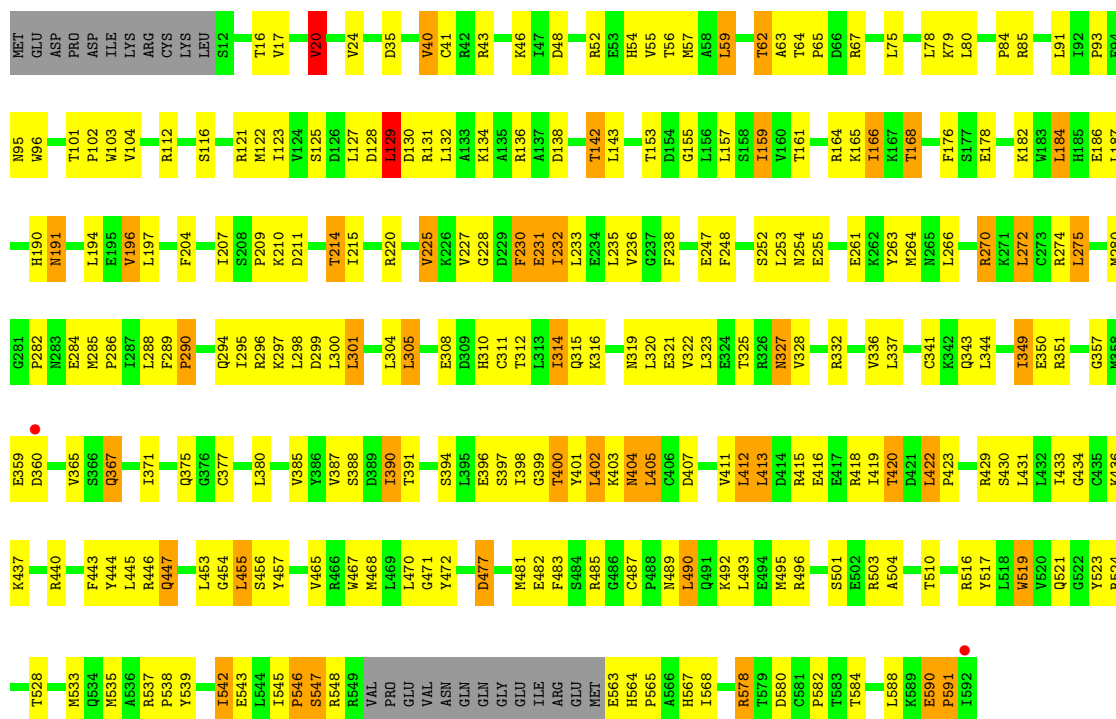
Chain B:





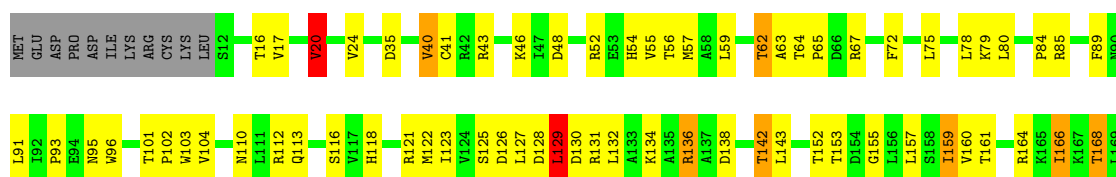
• Molecule 2: Coronatine-insensitiveprotein 1

Chain D:



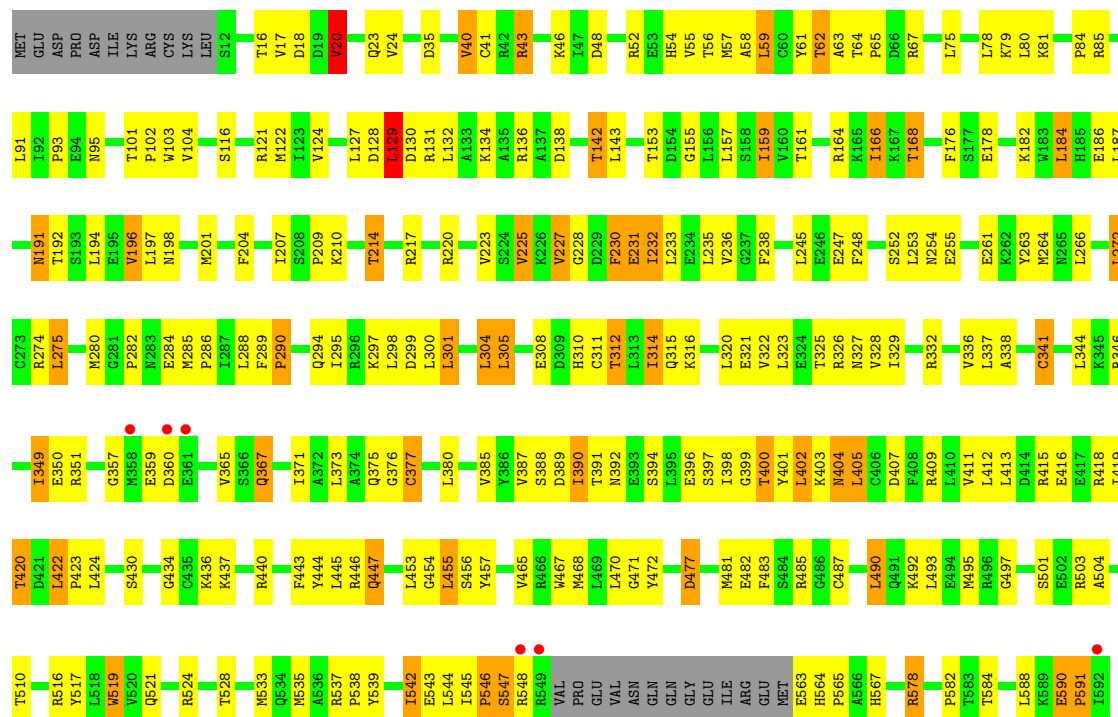
• Molecule 2: Coronatine-insensitiveprotein 1

Chain F:



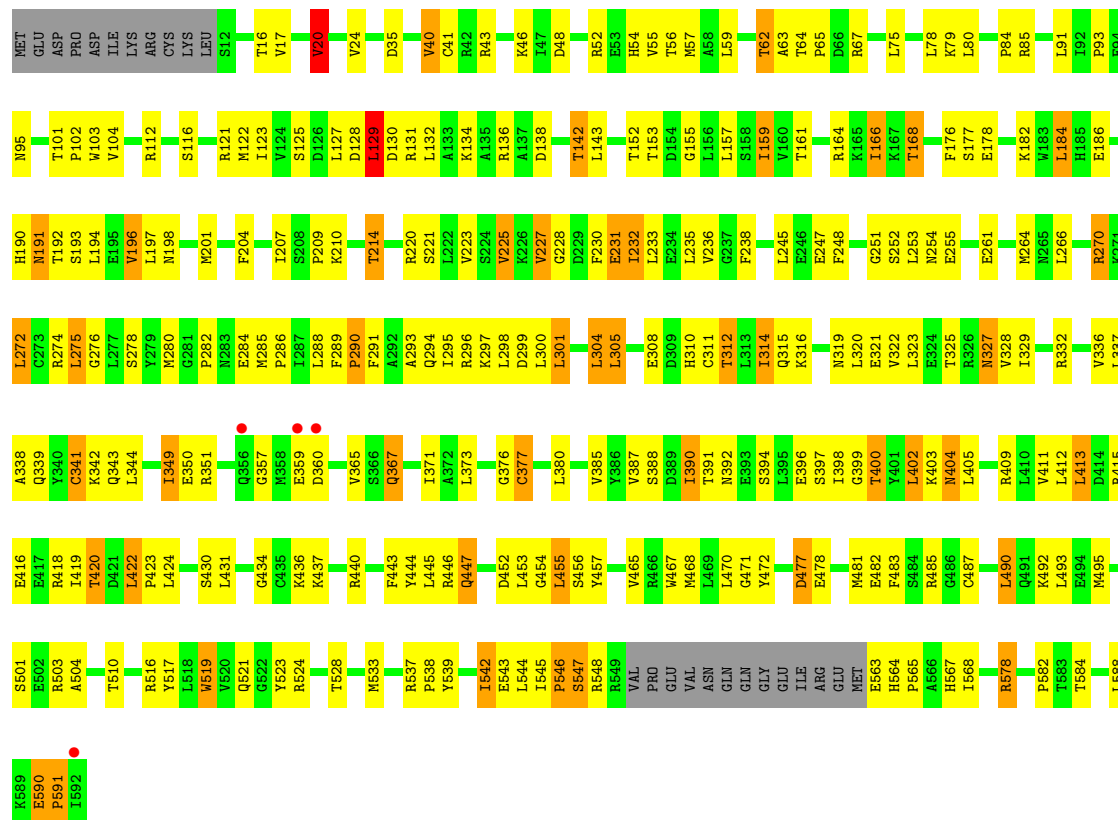


Chain J:



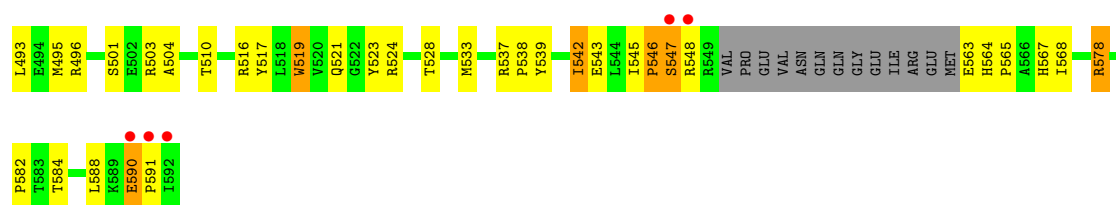
- Molecule 2: Coronatine-insensitive protein 1

Chain L:



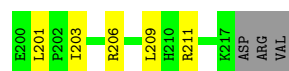
Chain N:





- Molecule 3: JAZ1 degron peptide

Chain Q:



- Molecule 3: JAZ1 degron peptide

Chain R:



- Molecule 3: JAZ1 degron peptide

Chain S:



- Molecule 3: JAZ1 degron peptide

Chain U:



- Molecule 3: JAZ1 degron peptide

Chain V:



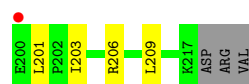
- Molecule 3: JAZ1 degron peptide

Chain W:



- Molecule 3: JAZ1 degron peptide

Chain X:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	123.17Å 220.76Å 149.54Å 90.00° 104.45° 90.00°	Depositor
Resolution (Å)	49.94 – 3.34 49.94 – 3.34	Depositor EDS
% Data completeness (in resolution range)	85.8 (49.94-3.34) 85.9 (49.94-3.34)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.96 (at 3.33Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.226 , 0.270 0.213 , 0.254	Depositor DCC
R_{free} test set	1871 reflections (1.95%)	DCC
Wilson B-factor (Å ²)	68.3	Xtriage
Anisotropy	0.387	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 23.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 103078 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	46877	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.02% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, OGK

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	0.44	0/1162	0.56	0/1571
1	C	0.45	0/1162	0.58	0/1571
1	E	0.45	0/1162	0.59	0/1571
1	G	0.52	0/1162	0.63	0/1571
1	I	0.44	0/1162	0.57	0/1571
1	K	0.43	0/1162	0.57	0/1571
1	M	0.50	0/1162	0.59	0/1571
1	O	0.45	0/1162	0.56	0/1571
2	B	0.49	0/4623	0.63	1/6238 (0.0%)
2	D	0.50	0/4623	0.64	1/6238 (0.0%)
2	F	0.46	0/4623	0.64	1/6238 (0.0%)
2	H	0.58	0/4566	0.83	2/6161 (0.0%)
2	J	0.49	0/4623	0.64	1/6238 (0.0%)
2	L	0.45	0/4623	0.63	1/6238 (0.0%)
2	N	0.58	0/4623	0.67	1/6238 (0.0%)
2	P	0.47	0/4623	0.63	1/6238 (0.0%)
3	Q	0.42	0/158	0.53	0/208
3	R	0.39	0/158	0.52	0/208
3	S	0.38	0/158	0.55	0/208
3	U	0.46	0/158	0.55	0/208
3	V	0.41	0/158	0.54	0/208
3	W	0.43	0/158	0.57	0/208
3	X	0.40	0/158	0.53	0/208
All	All	0.50	0/47329	0.65	9/63851 (0.0%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	129	LEU	CA-CB-CG	7.63	132.84	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	129	LEU	CA-CB-CG	7.51	132.58	115.30
2	L	129	LEU	CA-CB-CG	7.50	132.54	115.30
2	B	129	LEU	CA-CB-CG	7.40	132.32	115.30
2	N	129	LEU	CA-CB-CG	7.21	131.87	115.30
2	D	129	LEU	CA-CB-CG	7.03	131.47	115.30
2	P	129	LEU	CA-CB-CG	6.39	130.00	115.30
2	H	304	LEU	CA-CB-CG	-5.41	102.86	115.30
2	H	390	ILE	CB-CA-C	-5.06	101.47	111.60

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1146	0	1117	63	0
1	C	1146	0	1117	57	0
1	E	1146	0	1117	69	0
1	G	1146	0	1117	65	0
1	I	1146	0	1117	69	0
1	K	1146	0	1117	55	0
1	M	1146	0	1117	86	0
1	O	1146	0	1117	69	0
2	B	4541	0	4583	222	0
2	D	4541	0	4583	225	0
2	F	4541	0	4583	219	0
2	H	4486	0	4534	431	0
2	J	4541	0	4583	226	0
2	L	4541	0	4583	233	0
2	N	4541	0	4583	271	0
2	P	4541	0	4583	224	0
3	Q	156	0	171	7	0
3	R	156	0	171	5	0
3	S	156	0	171	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	U	156	0	171	6	0
3	V	156	0	171	5	0
3	W	156	0	171	7	0
3	X	156	0	171	5	0
4	B	23	0	27	7	0
4	D	23	0	27	7	0
4	F	23	0	27	7	0
4	H	23	0	27	16	0
4	J	23	0	27	7	0
4	L	23	0	27	7	0
4	N	23	0	27	6	0
4	P	23	0	27	8	0
5	B	20	0	0	3	0
5	D	20	0	0	2	0
5	F	20	0	0	2	0
5	H	20	0	0	3	0
5	J	20	0	0	2	0
5	L	20	0	0	2	0
5	N	20	0	0	3	0
5	P	20	0	0	3	0
All	All	46877	0	46964	2465	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 26.

All (2465) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:116:SER:HB2	2:H:142:THR:HG23	1.35	1.09
2:P:168:THR:HB	2:P:196:VAL:HG13	1.35	1.08
2:D:168:THR:HB	2:D:196:VAL:HG13	1.35	1.08
2:L:168:THR:HB	2:L:196:VAL:HG13	1.35	1.06
2:J:168:THR:HB	2:J:196:VAL:HG13	1.35	1.06
2:L:542:ILE:HD11	2:L:588:LEU:HD12	1.35	1.06
2:N:168:THR:HB	2:N:196:VAL:HG13	1.38	1.05
2:B:168:THR:HB	2:B:196:VAL:HG13	1.37	1.05
2:N:142:THR:HB	2:N:168:THR:HG23	1.38	1.05
1:M:93:ILE:HD12	1:M:97:THR:HG22	1.40	1.04
2:H:78:LEU:HD12	2:H:79:LYS:H	1.22	1.04
2:H:311:CYS:HB3	2:H:336:VAL:HG21	1.34	1.04
2:F:142:THR:HB	2:F:168:THR:HG23	1.40	1.04
2:J:542:ILE:HD11	2:J:588:LEU:HD12	1.40	1.04
2:D:542:ILE:HD11	2:D:588:LEU:HD12	1.40	1.04

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:93:ILE:HD12	1:A:97:THR:HG22	1.40	1.02
2:F:168:THR:HB	2:F:196:VAL:HG13	1.39	1.01
2:F:542:ILE:HD11	2:F:588:LEU:HD12	1.43	1.01
2:H:310:HIS:O	2:H:314:ILE:HG12	1.61	1.00
2:P:142:THR:HB	2:P:168:THR:HG23	1.43	1.00
2:N:542:ILE:HD11	2:N:588:LEU:HD12	1.40	0.99
2:P:542:ILE:HD11	2:P:588:LEU:HD12	1.45	0.98
2:F:93:PRO:HA	2:F:548:ARG:HB2	1.44	0.98
2:L:142:THR:HB	2:L:168:THR:HG23	1.46	0.98
2:P:357:GLY:HA2	2:P:415:ARG:HH22	1.29	0.97
2:L:93:PRO:HA	2:L:548:ARG:HB2	1.42	0.97
2:P:93:PRO:HA	2:P:548:ARG:HB2	1.44	0.97
1:M:159:PHE:O	1:M:160:GLU:HB2	1.63	0.96
2:F:357:GLY:HA2	2:F:415:ARG:HH22	1.30	0.96
2:H:392:ASN:O	2:H:396:GLU:HG2	1.66	0.96
2:B:357:GLY:HA2	2:B:415:ARG:HH22	1.31	0.96
1:K:93:ILE:HD12	1:K:97:THR:HG22	1.45	0.95
1:O:93:ILE:HD12	1:O:97:THR:HG22	1.49	0.95
2:H:152:THR:HG22	2:H:177:SER:HB2	1.47	0.95
2:D:93:PRO:HA	2:D:548:ARG:HB2	1.48	0.95
2:B:93:PRO:HA	2:B:548:ARG:HB2	1.48	0.94
2:J:93:PRO:HA	2:J:548:ARG:HB2	1.48	0.94
2:L:357:GLY:HA2	2:L:415:ARG:HH22	1.29	0.94
2:N:93:PRO:HA	2:N:548:ARG:HB2	1.50	0.94
1:K:159:PHE:O	1:K:160:GLU:HB2	1.66	0.94
2:J:357:GLY:HA2	2:J:415:ARG:HH22	1.30	0.94
1:E:93:ILE:HD12	1:E:97:THR:HG22	1.46	0.94
2:J:142:THR:HB	2:J:168:THR:HG23	1.49	0.93
2:H:95:ASN:H	2:H:95:ASN:HD22	1.09	0.93
2:B:142:THR:HB	2:B:168:THR:HG23	1.51	0.93
1:C:93:ILE:HD12	1:C:97:THR:HG22	1.49	0.93
1:I:93:ILE:HD12	1:I:97:THR:HG22	1.47	0.93
2:B:542:ILE:HD11	2:B:588:LEU:HD12	1.50	0.93
1:A:159:PHE:O	1:A:160:GLU:HB2	1.69	0.93
1:E:159:PHE:O	1:E:160:GLU:HB2	1.67	0.92
1:O:159:PHE:O	1:O:160:GLU:HB2	1.67	0.92
2:D:357:GLY:HA2	2:D:415:ARG:HH22	1.32	0.92
1:G:160:GLU:HG3	2:H:52:ARG:HH21	1.35	0.92
2:N:357:GLY:HA2	2:N:415:ARG:HH22	1.32	0.91
1:I:159:PHE:O	1:I:160:GLU:HB2	1.67	0.91
1:G:93:ILE:HD12	1:G:97:THR:HG22	1.52	0.91
2:H:286:PRO:HA	2:H:289:PHE:CE2	2.06	0.91

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:415:ARG:HH12	2:H:472:TYR:HE1	1.08	0.91
1:C:159:PHE:O	1:C:160:GLU:HB2	1.69	0.90
2:J:590:GLU:HB3	2:J:591:PRO:HD2	1.54	0.90
2:B:444:TYR:HA	2:B:471:GLY:HA3	1.54	0.90
2:B:210:LYS:O	2:B:214:THR:HG22	1.72	0.90
2:D:444:TYR:HA	2:D:471:GLY:HA3	1.53	0.89
2:F:590:GLU:HB3	2:F:591:PRO:HD2	1.54	0.89
2:L:590:GLU:HB3	2:L:591:PRO:HD2	1.54	0.89
2:D:590:GLU:HB3	2:D:591:PRO:HD2	1.55	0.89
2:J:419:ILE:HD13	2:J:446:ARG:HH12	1.38	0.88
2:N:444:TYR:HA	2:N:471:GLY:HA3	1.55	0.88
2:H:322:VAL:HG22	2:H:346:ARG:HB2	1.54	0.88
1:E:107:TYR:OH	2:L:294:GLN:NE2	2.07	0.88
2:P:590:GLU:HB3	2:P:591:PRO:HD2	1.53	0.88
2:N:590:GLU:HB3	2:N:591:PRO:HD2	1.54	0.87
2:B:590:GLU:HB3	2:B:591:PRO:HD2	1.54	0.87
2:F:444:TYR:HA	2:F:471:GLY:HA3	1.55	0.87
2:J:367:GLN:HB3	2:J:391:THR:HG22	1.58	0.86
2:L:444:TYR:HA	2:L:471:GLY:HA3	1.57	0.86
1:E:105:ALA:HB3	1:E:114:LEU:HD13	1.56	0.86
1:K:108:LEU:HD12	1:K:110:ILE:HD11	1.58	0.86
2:B:419:ILE:HD13	2:B:446:ARG:HH12	1.40	0.86
2:H:365:VAL:HG11	2:H:387:VAL:HG22	1.55	0.85
2:H:519:TRP:NE1	4:H:4100:OGK:H01	1.90	0.85
1:M:125:ILE:HG23	1:M:133:ILE:HD12	1.58	0.85
2:N:367:GLN:HB3	2:N:391:THR:HG22	1.57	0.85
2:J:444:TYR:HA	2:J:471:GLY:HA3	1.56	0.85
2:J:210:LYS:O	2:J:214:THR:HG22	1.77	0.85
2:L:349:ILE:HG13	2:L:385:VAL:HG22	1.57	0.85
2:P:286:PRO:HA	2:P:289:PHE:CE2	2.12	0.85
1:E:43:PRO:HG2	2:L:319:ASN:ND2	1.92	0.85
2:H:152:THR:CG2	2:H:177:SER:HB2	2.07	0.85
1:M:102:ILE:HD12	2:N:20:VAL:HG21	1.58	0.84
2:H:40:VAL:O	2:H:41:CYS:HB3	1.78	0.84
2:H:285:MET:HG2	2:H:286:PRO:HD3	1.57	0.84
2:H:415:ARG:NH1	2:H:472:TYR:CE1	2.46	0.84
2:H:440:ARG:HB3	2:H:467:TRP:HD1	1.43	0.84
1:G:153:ARG:HG3	1:G:157:TRP:CZ3	2.13	0.84
2:H:95:ASN:ND2	2:H:95:ASN:H	1.75	0.83
2:D:142:THR:HB	2:D:168:THR:HG23	1.58	0.83
2:B:116:SER:HB2	2:B:142:THR:HG23	1.59	0.83
2:H:199:PHE:CZ	2:H:227:VAL:HG23	2.13	0.83

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:125:ILE:HG23	1:I:133:ILE:HD12	1.58	0.83
2:N:328:VAL:HG22	2:N:359:GLU:HB2	1.61	0.82
2:D:210:LYS:O	2:D:214:THR:HG22	1.77	0.82
2:L:210:LYS:O	2:L:214:THR:HG22	1.80	0.82
2:F:210:LYS:O	2:F:214:THR:HG22	1.79	0.82
2:F:367:GLN:HB3	2:F:391:THR:HG22	1.61	0.82
2:L:367:GLN:HB3	2:L:391:THR:HG22	1.60	0.82
2:P:444:TYR:HA	2:P:471:GLY:HA3	1.61	0.82
2:D:116:SER:HB2	2:D:142:THR:HG23	1.62	0.82
1:I:108:LEU:HD12	1:I:110:ILE:HD11	1.62	0.81
1:M:108:LEU:HD12	1:M:110:ILE:HD11	1.62	0.81
2:H:387:VAL:HG11	2:H:390:ILE:HD13	1.60	0.81
1:A:108:LEU:HD12	1:A:110:ILE:HD11	1.61	0.81
2:D:367:GLN:HB3	2:D:391:THR:HG22	1.60	0.81
2:P:419:ILE:HD13	2:P:446:ARG:HH12	1.44	0.81
2:H:256:ASP:OD2	2:H:259:MET:HG2	1.80	0.81
2:J:286:PRO:HA	2:J:289:PHE:CE2	2.16	0.81
2:N:419:ILE:HD13	2:N:446:ARG:HH12	1.44	0.81
2:H:253:LEU:HD11	2:H:277:LEU:HD13	1.62	0.81
2:N:116:SER:HB2	2:N:142:THR:HG23	1.60	0.81
1:M:102:ILE:HG21	2:N:20:VAL:HG22	1.62	0.81
2:J:164:ARG:HE	2:N:112:ARG:CG	1.94	0.81
4:N:7100:OGK:H18A	4:N:7100:OGK:HN08	1.45	0.81
2:P:328:VAL:HG22	2:P:359:GLU:HB2	1.62	0.81
2:N:210:LYS:O	2:N:214:THR:HG22	1.81	0.80
2:D:419:ILE:HD13	2:D:446:ARG:HH12	1.44	0.80
2:P:210:LYS:O	2:P:214:THR:HG22	1.79	0.80
2:H:286:PRO:HA	2:H:289:PHE:CD2	2.15	0.80
2:L:286:PRO:HA	2:L:289:PHE:CE2	2.15	0.80
2:P:367:GLN:HB3	2:P:391:THR:HG22	1.63	0.80
2:H:361:GLU:HA	2:H:388:SER:OG	1.81	0.80
2:B:328:VAL:HG22	2:B:359:GLU:HB2	1.64	0.80
2:F:419:ILE:HD13	2:F:446:ARG:HH12	1.47	0.79
2:N:349:ILE:HG13	2:N:385:VAL:HG22	1.62	0.79
1:O:125:ILE:HG23	1:O:133:ILE:HD12	1.63	0.79
2:H:441:PHE:O	2:H:468:MET:HA	1.82	0.79
1:A:125:ILE:HG23	1:A:133:ILE:HD12	1.65	0.79
2:L:289:PHE:CD1	2:L:316:LYS:HD2	2.18	0.79
2:D:328:VAL:HG22	2:D:359:GLU:HB2	1.64	0.79
2:F:328:VAL:HG22	2:F:359:GLU:HB2	1.63	0.79
2:F:112:ARG:HG2	2:L:164:ARG:HD2	1.65	0.79
2:H:364:LEU:HD13	2:H:388:SER:OG	1.81	0.79

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:367:GLN:HB3	2:B:391:THR:HG22	1.63	0.79
2:F:116:SER:HB2	2:F:142:THR:HG23	1.65	0.78
4:B:1100:OGK:HN08	4:B:1100:OGK:H18A	1.48	0.78
2:B:286:PRO:HA	2:B:289:PHE:CE2	2.18	0.78
2:P:116:SER:HB2	2:P:142:THR:HG23	1.65	0.78
2:B:289:PHE:CD1	2:B:316:LYS:HD2	2.18	0.78
2:L:161:THR:HG22	2:L:186:GLU:HG2	1.65	0.78
2:J:328:VAL:HG22	2:J:359:GLU:HB2	1.66	0.78
2:J:349:ILE:HG13	2:J:385:VAL:HG22	1.64	0.78
2:H:283:ASN:O	2:H:286:PRO:HD2	1.83	0.78
1:M:96:ALA:HB1	2:N:14:VAL:HG12	1.65	0.78
2:L:419:ILE:HD13	2:L:446:ARG:HH12	1.49	0.78
2:H:411:VAL:HG13	2:H:444:TYR:HB3	1.66	0.77
2:D:349:ILE:HG13	2:D:385:VAL:HG22	1.63	0.77
2:L:116:SER:HB2	2:L:142:THR:HG23	1.64	0.77
2:H:519:TRP:HE1	4:H:4100:OGK:H01	1.46	0.77
2:D:289:PHE:CD1	2:D:316:LYS:HD2	2.19	0.77
2:F:286:PRO:HA	2:F:289:PHE:CE2	2.19	0.77
2:H:533:MET:CE	2:H:588:LEU:HD13	2.14	0.77
2:H:387:VAL:CG1	2:H:390:ILE:HD13	2.14	0.77
1:M:102:ILE:HG21	2:N:20:VAL:CG2	2.14	0.77
4:H:4100:OGK:H18A	4:H:4100:OGK:HN08	1.50	0.77
4:P:8100:OGK:HN08	4:P:8100:OGK:H18A	1.49	0.77
1:K:125:ILE:HG23	1:K:133:ILE:HD12	1.67	0.77
2:H:459:GLY:O	2:H:486:GLY:HA3	1.84	0.77
1:I:105:ALA:HB3	1:I:114:LEU:HD13	1.66	0.77
2:L:328:VAL:HG22	2:L:359:GLU:HB2	1.65	0.77
2:D:286:PRO:HA	2:D:289:PHE:CE2	2.20	0.77
2:N:492:LYS:HE3	2:N:517:TYR:CE2	2.20	0.77
2:N:286:PRO:HA	2:N:289:PHE:CE2	2.19	0.77
2:H:450:LEU:HD11	2:H:454:GLY:HA3	1.65	0.76
2:N:289:PHE:CD1	2:N:316:LYS:HD2	2.19	0.76
1:C:125:ILE:HG23	1:C:133:ILE:HD12	1.65	0.76
2:H:542:ILE:HD11	2:H:588:LEU:HB2	1.67	0.76
2:P:286:PRO:HA	2:P:289:PHE:CD2	2.20	0.76
4:D:2100:OGK:HN08	4:D:2100:OGK:H18A	1.49	0.76
2:J:164:ARG:HE	2:N:112:ARG:HG3	1.51	0.76
2:P:168:THR:HB	2:P:196:VAL:CG1	2.13	0.75
2:P:289:PHE:CD1	2:P:316:LYS:HD2	2.20	0.75
4:F:3100:OGK:HN08	4:F:3100:OGK:H18A	1.49	0.75
2:J:116:SER:HB2	2:J:142:THR:HG23	1.65	0.75
2:N:168:THR:HB	2:N:196:VAL:CG1	2.16	0.75

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:J:308:GLU:HG3	2:J:332:ARG:HH22	1.52	0.75
1:C:108:LEU:HD12	1:C:110:ILE:HD11	1.69	0.75
1:E:125:ILE:HG23	1:E:133:ILE:HD12	1.69	0.75
1:E:108:LEU:HD12	1:E:110:ILE:HD11	1.68	0.75
2:H:321:GLU:HA	2:H:344:LEU:HA	1.69	0.75
2:H:365:VAL:HG21	2:H:387:VAL:HG13	1.68	0.75
2:H:467:TRP:HH2	2:H:494:GLU:OE1	1.70	0.75
2:N:391:THR:HG23	2:N:394:SER:H	1.52	0.75
2:F:289:PHE:CD1	2:F:316:LYS:HD2	2.22	0.74
2:L:308:GLU:HG3	2:L:332:ARG:HH22	1.51	0.74
2:H:170:LEU:HD23	2:H:170:LEU:O	1.86	0.74
2:H:415:ARG:NH1	2:H:472:TYR:HE1	1.84	0.74
1:K:124:MET:O	1:K:128:LYS:HE2	1.87	0.74
2:B:270:ARG:HB3	1:O:107:TYR:CD1	2.22	0.74
1:O:108:LEU:HD12	1:O:110:ILE:HD11	1.70	0.74
2:H:298:LEU:HD22	2:H:300:LEU:HG	1.70	0.74
2:J:289:PHE:CD1	2:J:316:LYS:HD2	2.22	0.74
1:M:124:MET:O	1:M:128:LYS:HE2	1.88	0.74
3:S:203:ILE:HD12	3:S:203:ILE:H	1.53	0.74
1:E:43:PRO:CG	2:L:319:ASN:ND2	2.50	0.74
1:M:129:THR:O	1:M:133:ILE:HG12	1.86	0.74
2:N:84:PRO:HB3	2:N:517:TYR:OH	1.87	0.73
2:F:472:TYR:OH	3:S:201:LEU:HB2	1.88	0.73
2:B:161:THR:HG22	2:B:186:GLU:HG2	1.68	0.73
2:J:286:PRO:HA	2:J:289:PHE:CD2	2.24	0.73
4:L:6100:OGK:H18A	4:L:6100:OGK:HN08	1.52	0.73
2:L:168:THR:HB	2:L:196:VAL:CG1	2.16	0.73
2:J:164:ARG:HD2	2:N:112:ARG:HG2	1.70	0.73
2:B:391:THR:HG23	2:B:394:SER:H	1.53	0.73
2:F:308:GLU:HG3	2:F:332:ARG:HH22	1.53	0.73
1:E:124:MET:O	1:E:128:LYS:HE2	1.89	0.73
2:H:136:ARG:NE	2:H:136:ARG:HA	2.04	0.73
2:P:391:THR:HG23	2:P:394:SER:H	1.54	0.73
2:H:120:ARG:NH2	5:H:1102:PO4:O3	2.20	0.73
2:H:412:LEU:HD11	2:H:445:LEU:HD22	1.69	0.72
2:J:391:THR:HG23	2:J:394:SER:H	1.54	0.72
4:J:5100:OGK:HN08	4:J:5100:OGK:H18A	1.54	0.72
2:N:308:GLU:HG3	2:N:332:ARG:HH22	1.54	0.72
1:K:129:THR:O	1:K:133:ILE:HG12	1.89	0.72
2:B:253:LEU:HD12	2:B:280:MET:HB2	1.72	0.72
2:D:440:ARG:HB3	2:D:467:TRP:HE3	1.54	0.72
2:H:467:TRP:CH2	2:H:494:GLU:OE1	2.42	0.72

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:76:ARG:HG2	2:H:76:ARG:HH11	1.55	0.72
2:J:168:THR:HB	2:J:196:VAL:CG1	2.17	0.72
2:L:286:PRO:HA	2:L:289:PHE:CD2	2.24	0.72
1:I:124:MET:O	1:I:128:LYS:HE2	1.89	0.72
2:B:349:ILE:HG13	2:B:385:VAL:HG22	1.71	0.72
2:P:308:GLU:HG3	2:P:332:ARG:HH22	1.53	0.72
2:H:391:THR:OG1	2:H:393:GLU:HB2	1.90	0.72
2:D:319:ASN:ND2	1:G:43:PRO:HG2	2.03	0.72
3:W:203:ILE:HD12	3:W:203:ILE:H	1.55	0.72
2:H:78:LEU:HD12	2:H:79:LYS:N	2.01	0.71
1:G:108:LEU:HD12	1:G:110:ILE:HD11	1.72	0.71
2:D:308:GLU:HG3	2:D:332:ARG:HH22	1.55	0.71
2:P:161:THR:HG22	2:P:186:GLU:HG2	1.71	0.71
2:D:391:THR:HG23	2:D:394:SER:H	1.54	0.71
1:M:99:PHE:HD2	2:N:15:ALA:O	1.73	0.71
2:B:157:LEU:O	2:B:161:THR:HG23	1.91	0.71
2:H:334:LEU:CD2	2:H:373:LEU:HD22	2.21	0.71
2:F:286:PRO:HA	2:F:289:PHE:CD2	2.25	0.71
2:B:308:GLU:HG3	2:B:332:ARG:HH22	1.56	0.71
1:E:129:THR:O	1:E:133:ILE:HG12	1.89	0.71
2:P:349:ILE:HG13	2:P:385:VAL:HG22	1.72	0.71
3:V:203:ILE:HD12	3:V:203:ILE:H	1.55	0.71
2:H:540:TRP:CZ2	2:H:570:ALA:HB1	2.25	0.71
2:D:492:LYS:HE3	2:D:517:TYR:CE2	2.26	0.70
2:H:211:ASP:O	2:H:215:ILE:HG13	1.91	0.70
1:E:105:ALA:CB	1:E:114:LEU:HD13	2.21	0.70
2:N:55:VAL:HG23	2:N:75:LEU:HD21	1.72	0.70
1:M:112:ASN:C	1:M:114:LEU:H	1.95	0.70
2:J:161:THR:HG22	2:J:186:GLU:HG2	1.73	0.70
3:U:203:ILE:H	3:U:203:ILE:HD12	1.56	0.70
2:L:391:THR:HG23	2:L:394:SER:H	1.56	0.70
2:H:76:ARG:HH11	2:H:76:ARG:CG	2.04	0.70
2:F:161:THR:HG22	2:F:186:GLU:HG2	1.72	0.70
2:F:85:ARG:NH2	4:F:3100:OGK:O07	2.23	0.70
2:H:116:SER:CB	2:H:142:THR:HG23	2.16	0.70
2:B:492:LYS:HE3	2:B:517:TYR:CE2	2.25	0.70
2:H:59:LEU:HD22	2:H:61:TYR:HB2	1.73	0.70
2:H:311:CYS:CB	2:H:336:VAL:HG21	2.19	0.70
2:P:55:VAL:HG23	2:P:75:LEU:HD21	1.73	0.70
2:H:370:LEU:N	2:H:370:LEU:HD23	2.06	0.70
2:H:364:LEU:HD22	2:H:388:SER:H	1.54	0.70
2:H:191:ASN:HD21	2:H:194:LEU:H	1.37	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:118:HIS:CD2	2:H:144:LYS:HD3	2.26	0.70
1:C:124:MET:O	1:C:128:LYS:HE2	1.92	0.70
2:F:101:THR:HG22	2:F:128:ASP:OD1	1.91	0.70
2:H:450:LEU:CD1	2:H:454:GLY:HA3	2.22	0.69
2:J:419:ILE:HD13	2:J:446:ARG:NH1	2.07	0.69
2:H:325:THR:O	2:H:349:ILE:HA	1.91	0.69
2:D:286:PRO:HA	2:D:289:PHE:CD2	2.28	0.69
2:P:191:ASN:HD21	2:P:194:LEU:H	1.40	0.69
1:I:35:ASP:HB3	2:N:243:ALA:HB1	1.73	0.69
2:L:143:LEU:HD23	2:L:159:ILE:HD13	1.75	0.69
1:M:10:SER:HB2	1:M:52:LEU:HD23	1.74	0.69
2:B:101:THR:HG22	2:B:128:ASP:OD1	1.92	0.69
2:F:295:ILE:HG21	2:F:298:LEU:HD13	1.74	0.69
2:B:350:GLU:HB3	3:Q:209:LEU:HD21	1.74	0.69
2:N:521:GLN:HG3	2:N:567:HIS:HD2	1.55	0.69
2:D:161:THR:HG22	2:D:186:GLU:HG2	1.75	0.69
2:F:168:THR:HB	2:F:196:VAL:CG1	2.18	0.69
2:B:419:ILE:HD13	2:B:446:ARG:NH1	2.08	0.69
2:F:349:ILE:HG13	2:F:385:VAL:HG22	1.73	0.69
2:H:367:GLN:O	2:H:371:ILE:HG13	1.93	0.69
2:H:176:PHE:CZ	2:H:204:PHE:CZ	2.80	0.69
2:F:297:LYS:HG3	2:F:322:VAL:HB	1.75	0.69
2:B:286:PRO:HA	2:B:289:PHE:CD2	2.27	0.69
2:F:157:LEU:O	2:F:161:THR:HG23	1.92	0.69
2:H:116:SER:HB2	2:H:142:THR:CG2	2.18	0.69
1:M:93:ILE:HD12	1:M:97:THR:CG2	2.21	0.69
2:H:133:ALA:HB2	2:H:159:ILE:HG22	1.75	0.69
2:P:492:LYS:HE3	2:P:517:TYR:CE2	2.28	0.69
3:X:203:ILE:HD12	3:X:203:ILE:H	1.58	0.69
3:R:203:ILE:HD12	3:R:203:ILE:H	1.56	0.69
2:F:391:THR:HG23	2:F:394:SER:H	1.55	0.68
2:N:161:THR:HG22	2:N:186:GLU:HG2	1.76	0.68
2:H:414:ASP:HB3	2:H:446:ARG:NH2	2.08	0.68
2:N:286:PRO:HA	2:N:289:PHE:CD2	2.28	0.68
2:H:278:SER:O	2:H:280:MET:N	2.24	0.68
2:J:164:ARG:NE	2:N:112:ARG:CG	2.55	0.68
1:K:10:SER:HB2	1:K:52:LEU:HD23	1.76	0.68
3:Q:203:ILE:HD12	3:Q:203:ILE:H	1.57	0.68
2:H:519:TRP:HH2	2:H:567:HIS:CE1	2.10	0.68
1:A:124:MET:O	1:A:128:LYS:HE2	1.94	0.68
2:N:519:TRP:HE1	4:N:7100:OGK:H01	1.57	0.68
2:L:253:LEU:HD12	2:L:280:MET:HB2	1.74	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:P:101:THR:HG22	2:P:128:ASP:OD1	1.93	0.68
2:B:168:THR:HB	2:B:196:VAL:CG1	2.18	0.68
2:J:164:ARG:CD	2:N:112:ARG:HG2	2.24	0.68
4:N:7100:OGK:C18	4:N:7100:OGK:HN08	2.07	0.68
2:L:468:MET:HE3	2:L:470:LEU:HD21	1.76	0.68
1:M:100:GLU:HG2	2:N:15:ALA:HB2	1.76	0.68
2:J:297:LYS:HG3	2:J:322:VAL:HB	1.75	0.68
1:O:129:THR:O	1:O:133:ILE:HG12	1.94	0.68
2:D:57:MET:HE3	2:D:62:THR:HG22	1.76	0.68
2:P:547:SER:HB3	2:P:564:HIS:CB	2.24	0.67
2:B:143:LEU:HD23	2:B:159:ILE:HD13	1.75	0.67
2:H:444:TYR:HA	2:H:471:GLY:HA3	1.76	0.67
1:I:129:THR:O	1:I:133:ILE:HG12	1.94	0.67
2:B:547:SER:HB3	2:B:564:HIS:CB	2.25	0.67
2:J:295:ILE:HG21	2:J:298:LEU:HD13	1.76	0.67
2:J:101:THR:HG22	2:J:128:ASP:OD1	1.94	0.67
2:L:492:LYS:HE3	2:L:517:TYR:CE2	2.29	0.67
2:H:81:LYS:HG2	2:H:120:ARG:HD3	1.77	0.67
2:J:547:SER:HB3	2:J:564:HIS:CB	2.24	0.67
2:B:84:PRO:HB3	2:B:517:TYR:OH	1.94	0.67
2:N:282:PRO:HA	2:N:285:MET:HE2	1.77	0.67
2:P:295:ILE:HG21	2:P:298:LEU:HD13	1.75	0.67
1:C:40:ASN:HB3	1:K:13:GLY:O	1.94	0.67
2:J:337:LEU:HD12	2:J:341:CYS:SG	2.34	0.67
2:P:157:LEU:O	2:P:161:THR:HG23	1.93	0.67
2:L:487:CYS:HB3	2:L:490:LEU:HB2	1.76	0.67
2:J:350:GLU:HB3	3:U:209:LEU:HD21	1.77	0.67
2:N:388:SER:O	2:N:416:GLU:HG3	1.95	0.67
2:D:297:LYS:HG3	2:D:322:VAL:HB	1.77	0.67
2:H:52:ARG:NH1	2:H:72:PHE:CZ	2.63	0.67
2:L:547:SER:HB3	2:L:564:HIS:CB	2.25	0.67
2:H:85:ARG:NH2	4:H:4100:OGK:O07	2.29	0.66
2:D:547:SER:HB3	2:D:564:HIS:CB	2.25	0.66
1:O:105:ALA:HB3	1:O:114:LEU:HD13	1.77	0.66
2:N:157:LEU:O	2:N:161:THR:HG23	1.96	0.66
2:D:295:ILE:HG21	2:D:298:LEU:HD13	1.78	0.66
2:N:297:LYS:HG3	2:N:322:VAL:HB	1.78	0.66
2:H:533:MET:HE1	2:H:588:LEU:HD13	1.77	0.66
2:H:59:LEU:CD2	2:H:61:TYR:HB2	2.25	0.66
2:J:440:ARG:HB3	2:J:467:TRP:HE3	1.59	0.66
1:M:52:LEU:O	1:M:56:ILE:HG13	1.96	0.66
2:P:253:LEU:HD12	2:P:280:MET:HB2	1.77	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:398:ILE:HG23	2:H:402:LEU:HD11	1.76	0.66
2:N:295:ILE:HG21	2:N:298:LEU:HD13	1.78	0.66
1:A:93:ILE:HD12	1:A:97:THR:CG2	2.22	0.66
2:P:419:ILE:HD13	2:P:446:ARG:NH1	2.11	0.66
2:N:101:THR:HG22	2:N:128:ASP:OD1	1.96	0.66
2:J:492:LYS:HE3	2:J:517:TYR:CE2	2.30	0.66
2:F:455:LEU:HD22	2:F:483:PHE:HB2	1.78	0.66
2:J:157:LEU:O	2:J:161:THR:HG23	1.95	0.66
2:L:472:TYR:OH	3:V:201:LEU:HB2	1.96	0.66
1:G:10:SER:HB2	1:G:52:LEU:HD23	1.77	0.66
2:H:546:PRO:HD2	2:H:584:THR:O	1.96	0.65
2:P:282:PRO:HA	2:P:285:MET:HE2	1.78	0.65
2:J:487:CYS:HB3	2:J:490:LEU:HB2	1.79	0.65
2:D:472:TYR:OH	3:R:201:LEU:HB2	1.96	0.65
2:D:282:PRO:HA	2:D:285:MET:HE2	1.79	0.65
1:I:105:ALA:HB3	1:I:114:LEU:CD1	2.26	0.65
2:H:492:LYS:NZ	2:H:516:ARG:HH11	1.93	0.65
2:B:487:CYS:HB3	2:B:490:LEU:HB2	1.77	0.65
2:B:590:GLU:HB3	2:B:591:PRO:CD	2.27	0.65
2:H:40:VAL:O	2:H:41:CYS:CB	2.43	0.65
2:J:138:ASP:OD2	2:J:164:ARG:HG3	1.96	0.65
2:B:85:ARG:NH2	4:B:1100:OGK:O07	2.28	0.65
2:N:57:MET:HE3	2:N:62:THR:HG22	1.79	0.65
2:N:547:SER:HB3	2:N:564:HIS:CB	2.26	0.65
2:H:375:GLN:HG2	2:H:401:TYR:CD1	2.31	0.65
2:N:129:LEU:HD21	2:N:155:GLY:HA3	1.78	0.65
2:P:487:CYS:HB3	2:P:490:LEU:HB2	1.77	0.65
2:H:304:LEU:O	2:H:304:LEU:HD13	1.96	0.65
2:H:95:ASN:HD22	2:H:95:ASN:N	1.78	0.65
1:E:37:CYS:SG	2:L:296:ARG:NH2	2.69	0.65
2:J:164:ARG:NE	2:N:112:ARG:HG2	2.11	0.65
2:N:143:LEU:HD23	2:N:159:ILE:HD13	1.78	0.65
2:J:191:ASN:HD21	2:J:194:LEU:H	1.45	0.65
2:F:84:PRO:HB3	2:F:517:TYR:OH	1.96	0.65
1:C:10:SER:HB2	1:C:52:LEU:HD23	1.79	0.65
2:H:533:MET:HE3	2:H:588:LEU:HD13	1.78	0.65
2:N:487:CYS:HB3	2:N:490:LEU:HB2	1.77	0.65
2:F:350:GLU:HB3	3:S:209:LEU:HD21	1.77	0.65
2:H:545:ILE:HB	2:H:567:HIS:HB2	1.79	0.64
2:B:440:ARG:HB3	2:B:467:TRP:HE3	1.61	0.64
2:H:492:LYS:NZ	2:H:516:ARG:NH1	2.45	0.64
2:L:297:LYS:HG3	2:L:322:VAL:HB	1.80	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:N:399:GLY:O	2:N:434:GLY:HA3	1.97	0.64
2:H:52:ARG:NH1	2:H:72:PHE:CE2	2.66	0.64
2:H:442:ALA:HB1	2:H:469:LEU:HB3	1.78	0.64
2:P:440:ARG:HB3	2:P:467:TRP:HE3	1.62	0.64
1:C:113:LEU:O	1:C:117:THR:HG23	1.96	0.64
2:B:295:ILE:HG21	2:B:298:LEU:HD13	1.79	0.64
1:I:26:SER:OG	1:I:108:LEU:HB3	1.98	0.64
2:F:101:THR:OG1	2:F:102:PRO:HD3	1.96	0.64
2:D:101:THR:HG22	2:D:128:ASP:OD1	1.98	0.64
2:H:371:ILE:O	2:H:375:GLN:HG3	1.97	0.64
2:D:191:ASN:HD21	2:D:194:LEU:H	1.44	0.64
2:H:286:PRO:C	2:H:288:LEU:H	2.00	0.64
2:F:143:LEU:HD23	2:F:159:ILE:HD13	1.78	0.64
2:L:440:ARG:HB3	2:L:467:TRP:HE3	1.61	0.64
1:O:10:SER:HB2	1:O:52:LEU:HD23	1.79	0.64
2:J:253:LEU:HD12	2:J:280:MET:HB2	1.80	0.64
1:M:102:ILE:CD1	2:N:20:VAL:HG21	2.26	0.64
2:D:487:CYS:HB3	2:D:490:LEU:HB2	1.80	0.64
2:H:384:ALA:HB2	2:H:409:ARG:HG3	1.79	0.64
2:H:409:ARG:HB2	4:H:4100:OGK:H16B	1.79	0.64
2:P:85:ARG:NH2	4:P:8100:OGK:O07	2.29	0.64
2:F:40:VAL:O	2:F:41:CYS:HB3	1.96	0.64
2:H:104:VAL:HG21	2:H:128:ASP:HB2	1.80	0.64
2:F:253:LEU:HD12	2:F:280:MET:HB2	1.79	0.64
2:H:364:LEU:O	2:H:365:VAL:HG13	1.97	0.64
2:L:157:LEU:O	2:L:161:THR:HG23	1.98	0.64
1:A:10:SER:HB2	1:A:52:LEU:HD23	1.78	0.63
1:I:10:SER:HB2	1:I:52:LEU:HD23	1.78	0.63
2:B:521:GLN:HG3	2:B:567:HIS:HD2	1.63	0.63
1:A:12:ASP:OD2	1:A:49:SER:HB2	1.98	0.63
2:H:225:VAL:HG22	2:H:245:LEU:HD11	1.80	0.63
2:P:57:MET:HE3	2:P:62:THR:HG22	1.81	0.63
2:D:388:SER:O	2:D:416:GLU:HG3	1.99	0.63
2:N:521:GLN:HG3	2:N:567:HIS:CD2	2.33	0.63
2:N:298:LEU:HD23	2:N:300:LEU:HD11	1.81	0.63
2:F:57:MET:HE3	2:F:62:THR:HG22	1.80	0.63
1:O:124:MET:O	1:O:128:LYS:HE2	1.97	0.63
1:C:129:THR:O	1:C:133:ILE:HG12	1.99	0.63
2:L:191:ASN:HD21	2:L:194:LEU:H	1.46	0.63
1:A:129:THR:O	1:A:133:ILE:HG12	1.98	0.63
1:M:96:ALA:HB1	2:N:14:VAL:CG1	2.28	0.63
1:M:102:ILE:CG2	2:N:20:VAL:CG2	2.76	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:J:55:VAL:HG23	2:J:75:LEU:HD21	1.79	0.63
2:D:168:THR:HB	2:D:196:VAL:CG1	2.21	0.63
2:N:519:TRP:NE1	4:N:7100:OGK:H01	2.12	0.63
2:F:521:GLN:HG3	2:F:567:HIS:HD2	1.64	0.63
1:M:12:ASP:OD2	1:M:49:SER:HB2	1.99	0.63
2:P:311:CYS:HB3	2:P:336:VAL:HG21	1.80	0.63
1:K:102:ILE:HG12	1:K:117:THR:HB	1.80	0.63
2:H:357:GLY:C	2:H:359:GLU:H	2.02	0.63
2:N:419:ILE:HD13	2:N:446:ARG:NH1	2.13	0.63
2:D:294:GLN:NE2	1:G:107:TYR:OH	2.32	0.63
2:N:311:CYS:HB3	2:N:336:VAL:HG21	1.79	0.63
1:I:102:ILE:HG23	1:I:118:CYS:SG	2.39	0.63
2:J:590:GLU:HB3	2:J:591:PRO:CD	2.27	0.62
2:H:501:SER:HB3	2:H:524:ARG:NH1	2.14	0.62
2:H:230:PHE:HB3	2:H:235:LEU:HD21	1.81	0.62
2:F:547:SER:HB3	2:F:564:HIS:CB	2.28	0.62
2:B:191:ASN:HD21	2:B:194:LEU:H	1.47	0.62
2:D:590:GLU:HB3	2:D:591:PRO:CD	2.28	0.62
1:E:10:SER:HB2	1:E:52:LEU:HD23	1.79	0.62
1:E:45:PRO:HB2	2:L:291:PHE:HA	1.81	0.62
2:N:85:ARG:NH2	4:N:7100:OGK:O07	2.28	0.62
2:F:282:PRO:HA	2:F:285:MET:HE2	1.81	0.62
2:F:492:LYS:HE3	2:F:517:TYR:CE2	2.34	0.62
2:H:366:SER:OG	2:H:368:ARG:HG2	1.99	0.62
2:J:164:ARG:HE	2:N:112:ARG:HG2	1.64	0.62
2:D:190:HIS:CG	2:H:112:ARG:HH11	2.17	0.62
2:N:310:HIS:O	2:N:314:ILE:HG12	2.00	0.62
2:F:590:GLU:HB3	2:F:591:PRO:CD	2.28	0.62
2:L:138:ASP:OD2	2:L:164:ARG:HG3	1.98	0.62
2:L:84:PRO:HB3	2:L:517:TYR:OH	1.99	0.62
2:L:295:ILE:HG21	2:L:298:LEU:HD13	1.79	0.62
2:L:455:LEU:HD22	2:L:483:PHE:HB2	1.81	0.62
2:N:63:ALA:HB1	2:N:67:ARG:HD2	1.82	0.62
2:L:55:VAL:HG23	2:L:75:LEU:HD21	1.81	0.62
2:N:253:LEU:HD12	2:N:280:MET:HB2	1.81	0.62
2:J:521:GLN:HG3	2:J:567:HIS:HD2	1.65	0.62
2:H:544:LEU:O	2:H:546:PRO:HD3	2.00	0.62
2:F:468:MET:HE3	2:F:470:LEU:HD21	1.82	0.62
2:D:350:GLU:HB3	3:R:209:LEU:HD21	1.81	0.62
2:F:440:ARG:HB3	2:F:467:TRP:HE3	1.65	0.62
2:B:388:SER:O	2:B:416:GLU:HG3	2.00	0.62
2:F:487:CYS:HB3	2:F:490:LEU:HB2	1.80	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:419:ILE:HD13	2:D:446:ARG:NH1	2.13	0.62
2:P:129:LEU:HD21	2:P:155:GLY:HA3	1.82	0.61
2:H:101:THR:HG23	2:H:131:ARG:HH12	1.65	0.61
2:H:442:ALA:CB	2:H:469:LEU:HB3	2.30	0.61
2:P:184:LEU:HD12	2:P:207:ILE:HB	1.81	0.61
2:H:253:LEU:HD13	2:H:284:GLU:HG3	1.82	0.61
2:H:440:ARG:HB3	2:H:467:TRP:CD1	2.31	0.61
2:J:164:ARG:NE	2:N:112:ARG:HG3	2.15	0.61
2:N:455:LEU:HD22	2:N:483:PHE:HB2	1.82	0.61
2:F:399:GLY:O	2:F:434:GLY:HA3	2.00	0.61
2:F:365:VAL:HG23	2:F:387:VAL:HA	1.83	0.61
2:H:85:ARG:O	2:H:88:MET:HG2	2.01	0.61
2:B:282:PRO:HA	2:B:285:MET:HE2	1.82	0.61
1:K:52:LEU:O	1:K:56:ILE:HG13	2.01	0.61
2:N:101:THR:OG1	2:N:102:PRO:HD3	2.00	0.61
1:A:99:PHE:HZ	1:A:137:PHE:HE1	1.46	0.61
2:D:253:LEU:HD12	2:D:280:MET:HB2	1.81	0.61
2:H:89:PHE:CE1	4:H:4100:OGK:H17	2.35	0.61
2:L:419:ILE:HD13	2:L:446:ARG:NH1	2.15	0.61
1:M:137:PHE:CD1	2:N:17:VAL:HG21	2.36	0.61
2:J:57:MET:HE3	2:J:62:THR:HG22	1.82	0.61
2:L:184:LEU:HD12	2:L:207:ILE:HB	1.82	0.61
4:P:8100:OGK:C18	4:P:8100:OGK:HN08	2.13	0.61
1:E:48:THR:HG22	1:E:51:ILE:HB	1.82	0.61
2:P:310:HIS:O	2:P:314:ILE:HG12	2.00	0.61
2:H:461:TYR:C	2:H:463:PRO:HD3	2.20	0.61
1:O:137:PHE:CD1	2:P:17:VAL:HG21	2.35	0.61
2:P:533:MET:HE3	2:P:588:LEU:HD13	1.83	0.61
2:H:412:LEU:HD12	2:H:445:LEU:HA	1.83	0.61
1:M:26:SER:OG	1:M:108:LEU:HB3	2.00	0.61
2:F:419:ILE:HD13	2:F:446:ARG:NH1	2.16	0.61
2:D:227:VAL:CG1	2:D:228:GLY:N	2.64	0.61
2:H:40:VAL:HG23	2:H:40:VAL:O	1.99	0.61
2:L:85:ARG:NH2	4:L:6100:OGK:O07	2.30	0.61
2:P:590:GLU:HB3	2:P:591:PRO:CD	2.26	0.61
2:P:455:LEU:HD22	2:P:483:PHE:HB2	1.81	0.61
1:E:12:ASP:OD2	1:E:49:SER:HB2	2.00	0.61
2:J:129:LEU:HD21	2:J:155:GLY:HA3	1.82	0.61
2:L:590:GLU:HB3	2:L:591:PRO:CD	2.27	0.60
2:N:311:CYS:O	2:N:315:GLN:HB2	2.00	0.60
1:A:137:PHE:CD1	2:B:17:VAL:HG21	2.36	0.60
2:L:388:SER:O	2:L:416:GLU:HG3	2.01	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:542:ILE:CD1	2:L:588:LEU:HD12	2.21	0.60
2:J:101:THR:OG1	2:J:102:PRO:HD3	2.00	0.60
2:H:428:VAL:HG11	2:H:443:PHE:CZ	2.36	0.60
2:N:272:LEU:HD21	2:N:275:LEU:HB3	1.82	0.60
2:H:392:ASN:HD22	2:H:422:LEU:HD13	1.66	0.60
2:B:533:MET:HE3	2:B:588:LEU:HD13	1.82	0.60
1:E:52:LEU:O	1:E:56:ILE:HG13	2.01	0.60
2:H:124:VAL:O	2:H:151:PHE:HB3	2.02	0.60
1:E:6:ILE:HD12	1:E:23:ALA:HB3	1.82	0.60
2:L:227:VAL:CG1	2:L:228:GLY:N	2.64	0.60
2:N:367:GLN:O	2:N:371:ILE:HG22	2.01	0.60
2:J:547:SER:HB3	2:J:564:HIS:HB3	1.83	0.60
2:H:109:ASN:C	2:H:110:ASN:HD22	2.05	0.60
2:H:233:LEU:HA	2:H:236:VAL:HG23	1.82	0.60
2:P:388:SER:O	2:P:416:GLU:HG3	2.00	0.60
2:P:565:PRO:HG2	3:X:201:LEU:HD21	1.83	0.60
2:L:521:GLN:HG3	2:L:567:HIS:HD2	1.64	0.60
2:J:519:TRP:HE1	4:J:5100:OGK:H01	1.66	0.60
1:G:52:LEU:O	1:G:56:ILE:HG13	2.02	0.60
1:I:52:LEU:O	1:I:56:ILE:HG13	2.01	0.60
2:F:130:ASP:OD1	2:F:134:LYS:HE3	2.00	0.60
2:H:396:GLU:O	2:H:400:THR:HG23	2.01	0.60
2:D:91:LEU:O	2:D:567:HIS:HE1	1.83	0.60
1:G:137:PHE:HB3	2:H:17:VAL:HG21	1.82	0.60
2:B:311:CYS:O	2:B:315:GLN:HB2	2.01	0.60
2:H:425:ASP:OD2	2:H:451:THR:HG23	2.02	0.60
2:L:101:THR:HG22	2:L:128:ASP:OD1	2.01	0.60
2:B:365:VAL:HG23	2:B:387:VAL:HA	1.84	0.60
4:F:3100:OGK:HN08	4:F:3100:OGK:C18	2.15	0.60
2:D:84:PRO:HB3	2:D:517:TYR:OH	2.02	0.60
2:L:227:VAL:HG13	2:L:228:GLY:N	2.16	0.60
2:F:310:HIS:O	2:F:314:ILE:HG12	2.02	0.60
1:A:58:TYR:CD2	1:A:113:LEU:HD13	2.37	0.60
2:H:87:ALA:HB2	2:H:92:ILE:HG13	1.83	0.60
1:M:100:GLU:HG2	2:N:15:ALA:CB	2.31	0.60
2:D:101:THR:HG22	2:D:128:ASP:HB3	1.83	0.60
2:B:443:PHE:CE2	2:B:445:LEU:HD11	2.37	0.60
2:H:445:LEU:HD12	2:H:473:VAL:HG12	1.83	0.60
2:H:469:LEU:HD12	2:H:494:GLU:O	2.01	0.60
2:D:367:GLN:HB3	2:D:391:THR:CG2	2.31	0.60
2:P:547:SER:HB3	2:P:564:HIS:HB3	1.84	0.60
2:J:519:TRP:NE1	4:J:5100:OGK:H01	2.17	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:298:LEU:HD23	2:B:300:LEU:HD11	1.84	0.60
2:H:47:ILE:O	2:H:51:THR:HG23	2.02	0.60
2:H:76:ARG:NH1	2:H:76:ARG:CG	2.65	0.60
2:B:101:THR:OG1	2:B:102:PRO:HD3	2.02	0.60
1:G:30:ALA:O	1:G:33:VAL:HG22	2.01	0.60
2:B:468:MET:HE3	2:B:470:LEU:HD21	1.82	0.60
2:J:367:GLN:HB3	2:J:391:THR:CG2	2.31	0.59
1:G:153:ARG:HG2	1:G:153:ARG:NH1	2.17	0.59
1:M:99:PHE:HB2	2:N:15:ALA:CB	2.32	0.59
2:D:157:LEU:O	2:D:161:THR:HG23	2.02	0.59
1:M:158:ALA:HA	2:N:62:THR:HG21	1.84	0.59
1:E:30:ALA:O	1:E:33:VAL:HG22	2.02	0.59
2:B:297:LYS:HG3	2:B:322:VAL:HB	1.83	0.59
2:L:365:VAL:HG23	2:L:387:VAL:HA	1.84	0.59
1:M:48:THR:HG22	1:M:51:ILE:HB	1.84	0.59
2:N:91:LEU:O	2:N:567:HIS:HE1	1.85	0.59
2:N:275:LEU:HD11	2:N:288:LEU:HD21	1.83	0.59
2:H:22:GLU:HG2	2:H:47:ILE:CD1	2.32	0.59
2:B:121:ARG:NH2	5:B:1103:PO4:O4	2.34	0.59
2:N:65:PRO:HB3	2:N:103:TRP:CE3	2.37	0.59
2:H:230:PHE:CD1	2:H:235:LEU:HD21	2.37	0.59
2:J:388:SER:O	2:J:416:GLU:HG3	2.03	0.59
2:P:298:LEU:HD23	2:P:300:LEU:HD11	1.84	0.59
2:D:184:LEU:HD12	2:D:207:ILE:HB	1.84	0.59
2:D:40:VAL:O	2:D:41:CYS:HB3	2.02	0.59
2:H:392:ASN:O	2:H:396:GLU:CG	2.45	0.59
1:O:93:ILE:HD12	1:O:97:THR:CG2	2.29	0.59
2:D:101:THR:OG1	2:D:102:PRO:HD3	2.02	0.59
2:P:357:GLY:HA2	2:P:415:ARG:NH2	2.11	0.59
2:H:444:TYR:OH	2:H:496:ARG:HD3	2.01	0.59
2:N:367:GLN:HB3	2:N:391:THR:CG2	2.30	0.59
2:H:125:SER:HB2	2:H:128:ASP:H	1.67	0.59
2:L:129:LEU:HD21	2:L:155:GLY:HA3	1.85	0.59
1:G:130:PRO:HD3	2:H:36:SER:HB2	1.84	0.59
2:H:22:GLU:HG2	2:H:47:ILE:HD11	1.83	0.59
2:B:227:VAL:CG1	2:B:228:GLY:N	2.65	0.59
2:D:311:CYS:O	2:D:315:GLN:HB2	2.03	0.59
1:K:12:ASP:OD2	1:K:49:SER:HB2	2.01	0.59
1:A:52:LEU:O	1:A:56:ILE:HG13	2.01	0.59
1:E:158:ALA:HA	2:F:62:THR:HG21	1.84	0.59
2:J:351:ARG:O	2:J:351:ARG:HG3	2.03	0.59
2:P:468:MET:HE3	2:P:470:LEU:HD21	1.83	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:412:LEU:HD11	2:H:445:LEU:CD2	2.33	0.59
1:E:43:PRO:HG3	2:L:319:ASN:HD21	1.67	0.59
2:N:65:PRO:HA	2:N:103:TRP:CZ3	2.38	0.59
2:D:261:GLU:HG2	2:D:264:MET:HG3	1.85	0.59
2:L:197:LEU:O	2:L:225:VAL:HA	2.02	0.59
2:B:261:GLU:HG2	2:B:264:MET:HG3	1.85	0.59
2:P:91:LEU:O	2:P:567:HIS:HE1	1.86	0.58
2:J:357:GLY:HA2	2:J:415:ARG:NH2	2.12	0.58
1:I:93:ILE:HD12	1:I:97:THR:CG2	2.26	0.58
2:B:519:TRP:NE1	4:B:1100:OGK:H01	2.18	0.58
2:N:404:ASN:HA	2:N:437:LYS:HD2	1.85	0.58
1:C:12:ASP:OD2	1:C:49:SER:HB2	2.03	0.58
1:G:153:ARG:HG2	1:G:153:ARG:HH11	1.67	0.58
2:D:319:ASN:ND2	1:G:43:PRO:CG	2.65	0.58
2:L:547:SER:HB3	2:L:564:HIS:HB3	1.86	0.58
1:C:52:LEU:O	1:C:56:ILE:HG13	2.02	0.58
2:P:311:CYS:O	2:P:315:GLN:HB2	2.04	0.58
1:K:158:ALA:HA	2:L:62:THR:HG21	1.84	0.58
1:M:105:ALA:HB2	1:M:113:LEU:HD23	1.83	0.58
2:F:227:VAL:CG1	2:F:228:GLY:N	2.66	0.58
2:P:197:LEU:O	2:P:225:VAL:HA	2.03	0.58
1:E:137:PHE:CD1	2:F:17:VAL:HG21	2.39	0.58
2:B:57:MET:HE3	2:B:62:THR:HG22	1.85	0.58
2:J:233:LEU:O	2:J:236:VAL:HG23	2.04	0.58
2:P:337:LEU:HD12	2:P:341:CYS:SG	2.44	0.58
2:F:80:LEU:HB2	2:F:122:MET:CE	2.33	0.58
2:J:533:MET:HE3	2:J:588:LEU:HD13	1.85	0.58
2:F:357:GLY:HA2	2:F:415:ARG:NH2	2.12	0.58
2:L:357:GLY:HA2	2:L:415:ARG:NH2	2.11	0.58
2:P:84:PRO:HB3	2:P:517:TYR:OH	2.04	0.58
2:B:63:ALA:HB1	2:B:67:ARG:HD2	1.85	0.58
2:H:311:CYS:HB3	2:H:336:VAL:CG2	2.22	0.58
2:P:519:TRP:NE1	4:P:8100:OGK:H01	2.18	0.58
4:J:5100:OGK:HN08	4:J:5100:OGK:C18	2.16	0.58
2:J:91:LEU:O	2:J:567:HIS:HE1	1.87	0.58
2:B:227:VAL:HG13	2:B:228:GLY:N	2.19	0.58
1:G:98:LEU:HD21	1:G:120:THR:HG22	1.85	0.58
2:F:388:SER:O	2:F:416:GLU:HG3	2.03	0.58
2:H:390:ILE:HD11	2:H:410:LEU:HD11	1.83	0.58
2:N:80:LEU:HB2	2:N:122:MET:CE	2.34	0.58
1:O:12:ASP:OD2	1:O:49:SER:HB2	2.04	0.58
2:F:129:LEU:HD21	2:F:155:GLY:HA3	1.86	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:282:PRO:HA	2:L:285:MET:HE2	1.85	0.58
2:P:519:TRP:HE1	4:P:8100:OGK:H01	1.67	0.58
2:J:40:VAL:O	2:J:41:CYS:HB3	2.03	0.58
2:P:272:LEU:HD21	2:P:275:LEU:HB3	1.84	0.58
2:H:519:TRP:CH2	2:H:567:HIS:CE1	2.92	0.58
2:D:440:ARG:HB3	2:D:467:TRP:CE3	2.37	0.58
1:E:158:ALA:HA	2:F:62:THR:CG2	2.34	0.58
2:H:25:MET:HG2	2:H:47:ILE:HG22	1.86	0.58
1:G:12:ASP:OD2	1:G:49:SER:HB2	2.04	0.58
2:P:399:GLY:O	2:P:434:GLY:HA3	2.03	0.58
2:H:305:LEU:HD23	2:H:305:LEU:H	1.69	0.58
2:B:444:TYR:HA	2:B:471:GLY:CA	2.30	0.58
2:L:367:GLN:HB3	2:L:391:THR:CG2	2.32	0.58
2:D:547:SER:HB3	2:D:564:HIS:HB3	1.85	0.58
2:D:227:VAL:HG13	2:D:228:GLY:N	2.19	0.58
1:K:137:PHE:CD1	2:L:17:VAL:HG21	2.39	0.58
2:D:129:LEU:HD21	2:D:155:GLY:HA3	1.85	0.58
2:L:130:ASP:OD1	2:L:134:LYS:HE3	2.03	0.58
2:L:533:MET:HE3	2:L:588:LEU:HD13	1.85	0.57
2:F:446:ARG:HG2	2:F:447:GLN:H	1.69	0.57
1:I:134:ARG:HB2	1:I:139:ILE:O	2.04	0.57
2:B:129:LEU:HD21	2:B:155:GLY:HA3	1.85	0.57
2:B:477:ASP:OD2	2:B:504:ALA:HB2	2.04	0.57
2:H:38:SER:O	2:H:45:PHE:HB2	2.04	0.57
2:D:533:MET:HE3	2:D:588:LEU:HD13	1.87	0.57
2:P:101:THR:HG22	2:P:128:ASP:HB3	1.84	0.57
2:F:396:GLU:O	2:F:400:THR:HB	2.04	0.57
2:N:184:LEU:HD12	2:N:207:ILE:HB	1.86	0.57
1:O:91:MET:SD	1:O:117:THR:HG22	2.44	0.57
2:H:519:TRP:HE1	4:H:4100:OGK:C01	2.16	0.57
1:M:158:ALA:HA	2:N:62:THR:CG2	2.34	0.57
2:B:396:GLU:O	2:B:400:THR:HB	2.04	0.57
1:G:159:PHE:CE2	2:H:68:LEU:HD13	2.39	0.57
2:H:444:TYR:CD1	2:H:471:GLY:HA2	2.39	0.57
1:O:134:ARG:HB2	1:O:139:ILE:O	2.04	0.57
2:J:197:LEU:O	2:J:225:VAL:HA	2.04	0.57
2:B:91:LEU:O	2:B:567:HIS:HE1	1.87	0.57
2:J:184:LEU:HD12	2:J:207:ILE:HB	1.86	0.57
2:B:519:TRP:HE1	4:B:1100:OGK:H01	1.70	0.57
2:B:270:ARG:CB	1:O:107:TYR:CD1	2.88	0.57
2:B:101:THR:HG22	2:B:128:ASP:HB3	1.85	0.57
2:D:233:LEU:O	2:D:236:VAL:HG23	2.04	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:55:VAL:HG23	2:F:75:LEU:HD21	1.87	0.57
2:N:590:GLU:HB3	2:N:591:PRO:CD	2.27	0.57
1:I:105:ALA:CB	1:I:114:LEU:HD13	2.33	0.57
2:D:365:VAL:HG23	2:D:387:VAL:HA	1.85	0.57
2:P:297:LYS:HG3	2:P:322:VAL:HB	1.85	0.57
2:H:587:VAL:HG12	2:H:587:VAL:O	2.04	0.57
2:B:547:SER:HB3	2:B:564:HIS:HB3	1.86	0.57
1:M:134:ARG:HB2	1:M:139:ILE:O	2.04	0.57
2:H:395:LEU:HB3	2:H:427:GLY:O	2.04	0.57
2:L:93:PRO:HA	2:L:548:ARG:CB	2.26	0.57
2:J:101:THR:CG2	2:J:128:ASP:OD1	2.53	0.57
1:O:158:ALA:HA	2:P:62:THR:HG21	1.87	0.57
2:F:227:VAL:HG13	2:F:228:GLY:N	2.19	0.57
2:H:121:ARG:NH2	5:H:1103:PO4:O4	2.38	0.57
2:B:138:ASP:OD2	2:B:164:ARG:HG3	2.05	0.57
2:D:130:ASP:OD1	2:D:134:LYS:HE3	2.04	0.57
2:J:344:LEU:HD23	2:J:380:LEU:HD21	1.87	0.57
2:N:191:ASN:HD21	2:N:194:LEU:H	1.50	0.57
1:O:48:THR:HG22	1:O:51:ILE:HB	1.87	0.57
2:H:95:ASN:ND2	2:H:95:ASN:N	2.44	0.57
2:J:84:PRO:HB3	2:J:517:TYR:OH	2.05	0.57
2:P:40:VAL:O	2:P:41:CYS:HB3	2.02	0.57
2:N:261:GLU:HG2	2:N:264:MET:HG3	1.87	0.57
1:I:160:GLU:HA	1:I:160:GLU:OE2	2.05	0.56
2:H:407:ASP:OD1	2:H:440:ARG:HD2	2.05	0.56
2:H:542:ILE:CD1	2:H:588:LEU:HB2	2.35	0.56
1:E:134:ARG:HB2	1:E:139:ILE:O	2.05	0.56
1:K:6:ILE:HD12	1:K:23:ALA:HB3	1.86	0.56
2:P:396:GLU:O	2:P:400:THR:HB	2.04	0.56
1:A:48:THR:HG22	1:A:51:ILE:HB	1.86	0.56
2:J:143:LEU:HD23	2:J:159:ILE:HD13	1.87	0.56
2:L:311:CYS:O	2:L:315:GLN:HB2	2.04	0.56
1:E:160:GLU:HA	1:E:160:GLU:OE2	2.05	0.56
1:M:102:ILE:HD12	2:N:20:VAL:CG2	2.31	0.56
2:F:91:LEU:O	2:F:567:HIS:HE1	1.88	0.56
2:J:101:THR:HG22	2:J:128:ASP:HB3	1.86	0.56
2:H:225:VAL:HG21	2:H:238:PHE:HZ	1.70	0.56
2:D:396:GLU:O	2:D:400:THR:HB	2.05	0.56
2:L:357:GLY:CA	2:L:415:ARG:HH22	2.13	0.56
2:H:365:VAL:CG2	2:H:387:VAL:HG13	2.34	0.56
1:G:130:PRO:O	1:G:134:ARG:HG3	2.05	0.56
2:P:521:GLN:HG3	2:P:567:HIS:HD2	1.69	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:519:TRP:HE1	4:D:2100:OGK:H01	1.69	0.56
2:L:519:TRP:NE1	4:L:6100:OGK:H01	2.20	0.56
2:N:547:SER:HB3	2:N:564:HIS:HB3	1.86	0.56
1:O:52:LEU:O	1:O:56:ILE:HG13	2.05	0.56
2:F:443:PHE:CE2	2:F:445:LEU:HD11	2.40	0.56
2:F:261:GLU:HG2	2:F:264:MET:HG3	1.87	0.56
2:B:233:LEU:O	2:B:236:VAL:HG23	2.04	0.56
2:N:223:VAL:O	2:N:245:LEU:HD12	2.06	0.56
2:F:311:CYS:O	2:F:315:GLN:HB2	2.06	0.56
1:K:30:ALA:O	1:K:33:VAL:HG22	2.05	0.56
2:J:565:PRO:HG2	3:U:201:LEU:HD21	1.86	0.56
2:P:350:GLU:HB3	3:X:209:LEU:HD21	1.86	0.56
2:J:272:LEU:HD21	2:J:275:LEU:HB3	1.88	0.56
1:M:58:TYR:CD2	1:M:113:LEU:HD13	2.40	0.56
2:P:54:HIS:HE1	2:P:56:THR:OG1	1.88	0.56
2:P:227:VAL:CG1	2:P:228:GLY:N	2.68	0.56
2:H:411:VAL:CG1	2:H:444:TYR:HB3	2.35	0.56
1:E:43:PRO:HG2	2:L:319:ASN:HD22	1.68	0.56
2:H:82:GLY:HA3	2:H:122:MET:HG2	1.88	0.56
2:F:101:THR:CG2	2:F:128:ASP:OD1	2.54	0.56
2:D:298:LEU:HD23	2:D:300:LEU:HD11	1.87	0.56
2:L:311:CYS:HB3	2:L:336:VAL:HG21	1.88	0.56
2:D:121:ARG:NH2	5:D:1103:PO4:O4	2.37	0.56
2:H:276:GLY:HA3	2:H:299:ASP:HB3	1.87	0.56
2:P:93:PRO:HA	2:P:548:ARG:CB	2.29	0.56
2:D:357:GLY:HA2	2:D:415:ARG:NH2	2.14	0.56
2:H:334:LEU:HD22	2:H:373:LEU:HD22	1.85	0.56
2:N:40:VAL:O	2:N:41:CYS:HB3	2.04	0.56
1:C:48:THR:HG22	1:C:51:ILE:HB	1.86	0.56
2:F:482:GLU:O	2:F:485:ARG:HG2	2.05	0.56
2:N:396:GLU:O	2:N:400:THR:HB	2.05	0.56
2:D:404:ASN:HA	2:D:437:LYS:HD2	1.86	0.56
1:I:30:ALA:O	1:I:33:VAL:HG22	2.06	0.56
2:P:404:ASN:HA	2:P:437:LYS:HD2	1.88	0.56
2:N:477:ASP:OD2	2:N:504:ALA:HB2	2.05	0.56
2:B:472:TYR:OH	3:Q:201:LEU:HB2	2.05	0.56
2:H:447:GLN:HG3	2:H:447:GLN:O	2.05	0.56
2:D:444:TYR:HA	2:D:471:GLY:CA	2.31	0.56
2:H:384:ALA:CB	2:H:409:ARG:HG3	2.35	0.56
2:H:199:PHE:CE1	2:H:227:VAL:HG23	2.41	0.56
2:L:519:TRP:HE1	4:L:6100:OGK:H01	1.71	0.56
2:L:337:LEU:HD12	2:L:341:CYS:SG	2.46	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:P:443:PHE:CE2	2:P:445:LEU:HD11	2.40	0.56
2:H:326:ARG:HA	2:H:350:GLU:O	2.05	0.56
1:C:158:ALA:HA	2:D:62:THR:CG2	2.36	0.56
2:J:365:VAL:HG23	2:J:387:VAL:HA	1.87	0.56
2:F:275:LEU:HD11	2:F:288:LEU:HD21	1.88	0.56
2:D:138:ASP:OD2	2:D:164:ARG:HG3	2.06	0.56
1:G:48:THR:HG22	1:G:51:ILE:HB	1.88	0.56
2:H:343:GLN:CD	2:H:343:GLN:H	2.09	0.56
2:H:285:MET:CG	2:H:286:PRO:HD3	2.34	0.56
2:H:469:LEU:HD21	4:H:4100:OGK:H05	1.88	0.56
2:P:367:GLN:O	2:P:371:ILE:HG22	2.05	0.56
2:N:321:GLU:HA	2:N:344:LEU:HA	1.87	0.56
2:N:344:LEU:HD23	2:N:380:LEU:HD21	1.86	0.56
2:F:191:ASN:HD21	2:F:194:LEU:H	1.53	0.56
2:N:468:MET:HE3	2:N:470:LEU:HD21	1.88	0.56
2:B:40:VAL:O	2:B:41:CYS:HB3	2.06	0.56
2:L:396:GLU:O	2:L:400:THR:HB	2.06	0.56
2:F:93:PRO:HA	2:F:548:ARG:CB	2.29	0.56
2:D:411:VAL:HG22	2:D:444:TYR:HB3	1.87	0.56
1:M:96:ALA:CB	2:N:14:VAL:HG12	2.34	0.56
2:F:519:TRP:NE1	4:F:3100:OGK:H01	2.21	0.56
1:A:95:GLN:HE21	1:A:124:MET:HE3	1.71	0.56
2:H:462:SER:N	2:H:463:PRO:HD3	2.20	0.56
1:K:158:ALA:HA	2:L:62:THR:CG2	2.35	0.56
1:A:134:ARG:HB2	1:A:139:ILE:O	2.06	0.56
2:B:310:HIS:O	2:B:314:ILE:HG12	2.05	0.56
1:I:12:ASP:OD2	1:I:49:SER:HB2	2.05	0.56
2:F:184:LEU:HD12	2:F:207:ILE:HB	1.88	0.56
2:H:291:PHE:H	2:H:291:PHE:HD2	1.53	0.56
2:H:390:ILE:HD12	2:H:410:LEU:HD21	1.88	0.55
2:N:546:PRO:HD3	2:N:584:THR:O	2.06	0.55
2:B:55:VAL:HG23	2:B:75:LEU:HD21	1.88	0.55
2:P:351:ARG:HG3	2:P:351:ARG:O	2.06	0.55
2:L:272:LEU:HD21	2:L:275:LEU:HB3	1.87	0.55
1:G:125:ILE:HD11	2:H:44:TRP:HH2	1.71	0.55
2:L:399:GLY:O	2:L:434:GLY:HA3	2.06	0.55
1:C:30:ALA:O	1:C:33:VAL:HG22	2.05	0.55
2:L:411:VAL:HG22	2:L:444:TYR:HB3	1.88	0.55
2:D:519:TRP:NE1	4:D:2100:OGK:H01	2.20	0.55
2:H:76:ARG:O	2:H:114:LEU:HD12	2.07	0.55
2:B:397:SER:O	2:B:400:THR:HG22	2.06	0.55
2:P:397:SER:O	2:P:400:THR:HG22	2.06	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:40:VAL:O	2:L:41:CYS:HB3	2.06	0.55
2:J:396:GLU:O	2:J:400:THR:HB	2.05	0.55
1:E:45:PRO:HG3	2:L:293:ALA:HB3	1.87	0.55
1:G:153:ARG:HH11	1:G:153:ARG:CG	2.19	0.55
2:J:282:PRO:HA	2:J:285:MET:HE2	1.88	0.55
4:N:7100:OGK:C18	4:N:7100:OGK:N08	2.69	0.55
4:L:6100:OGK:HN08	4:L:6100:OGK:C18	2.17	0.55
2:H:477:ASP:O	2:H:481:MET:HG2	2.06	0.55
2:H:366:SER:HB3	2:H:368:ARG:HG3	1.88	0.55
1:O:153:ARG:NH2	2:P:539:TYR:CE1	2.74	0.55
2:F:519:TRP:HE1	4:F:3100:OGK:H01	1.71	0.55
1:G:26:SER:OG	1:G:108:LEU:HB3	2.06	0.55
2:J:261:GLU:HG2	2:J:264:MET:HG3	1.89	0.55
1:E:45:PRO:CD	2:L:294:GLN:HB2	2.37	0.55
2:L:101:THR:HG22	2:L:128:ASP:HB3	1.88	0.55
1:A:158:ALA:HA	2:B:62:THR:HG21	1.88	0.55
2:L:398:ILE:HG23	2:L:402:LEU:HD11	1.89	0.55
1:I:6:ILE:HD12	1:I:23:ALA:HB3	1.88	0.55
2:J:446:ARG:HG2	2:J:447:GLN:H	1.71	0.55
4:D:2100:OGK:C18	4:D:2100:OGK:HN08	2.19	0.55
2:P:101:THR:CG2	2:P:128:ASP:OD1	2.54	0.55
2:F:310:HIS:O	2:F:314:ILE:CG1	2.55	0.55
1:K:48:THR:HG22	1:K:51:ILE:HB	1.88	0.55
2:J:95:ASN:O	2:J:582:PRO:HG3	2.06	0.55
2:D:143:LEU:HD23	2:D:159:ILE:HD13	1.89	0.55
1:A:6:ILE:HD12	1:A:23:ALA:HB3	1.87	0.55
2:B:296:ARG:NH2	1:O:37:CYS:SG	2.80	0.55
2:H:232:ILE:HG13	2:H:253:LEU:HD21	1.89	0.55
2:F:367:GLN:HB3	2:F:391:THR:CG2	2.34	0.55
1:O:105:ALA:CB	1:O:114:LEU:HD13	2.37	0.55
2:H:374:ALA:HB2	2:H:398:ILE:HD12	1.88	0.55
1:C:102:ILE:HG12	1:C:117:THR:HB	1.89	0.55
2:P:310:HIS:O	2:P:314:ILE:CG1	2.55	0.55
1:A:30:ALA:O	1:A:33:VAL:HG22	2.06	0.55
2:J:311:CYS:O	2:J:315:GLN:HB2	2.06	0.55
2:P:261:GLU:HG2	2:P:264:MET:HG3	1.89	0.55
2:H:382:TYR:C	2:H:382:TYR:CD1	2.79	0.55
1:M:102:ILE:HB	2:N:20:VAL:HG21	1.89	0.55
1:O:48:THR:HG22	1:O:51:ILE:CG1	2.37	0.55
2:D:80:LEU:HB2	2:D:122:MET:CE	2.37	0.55
1:M:30:ALA:O	1:M:33:VAL:HG22	2.07	0.55
1:G:6:ILE:HD12	1:G:23:ALA:HB3	1.89	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:197:LEU:O	2:B:225:VAL:HA	2.07	0.55
1:A:91:MET:O	1:A:93:ILE:N	2.40	0.55
2:J:444:TYR:HA	2:J:471:GLY:CA	2.32	0.55
2:B:547:SER:HB3	2:B:564:HIS:CG	2.42	0.55
2:J:130:ASP:OD1	2:J:134:LYS:HE3	2.06	0.55
2:N:440:ARG:HB3	2:N:467:TRP:HE3	1.72	0.55
2:N:365:VAL:HG23	2:N:387:VAL:HA	1.88	0.55
2:H:232:ILE:HG13	2:H:253:LEU:CD2	2.37	0.55
2:D:319:ASN:HD22	1:G:43:PRO:HG2	1.71	0.55
2:J:298:LEU:HD23	2:J:300:LEU:HD11	1.88	0.55
2:B:311:CYS:HB3	2:B:336:VAL:HG21	1.89	0.55
2:L:57:MET:HE2	2:L:62:THR:HG22	1.88	0.55
2:N:397:SER:O	2:N:400:THR:HG22	2.07	0.55
1:G:129:THR:HG23	1:G:132:GLU:CD	2.27	0.55
2:D:197:LEU:O	2:D:225:VAL:HA	2.06	0.55
2:B:20:VAL:O	2:B:24:VAL:HG23	2.07	0.55
1:O:30:ALA:O	1:O:33:VAL:HG22	2.06	0.55
2:L:444:TYR:HA	2:L:471:GLY:CA	2.34	0.54
2:P:547:SER:HB3	2:P:564:HIS:CG	2.42	0.54
2:B:101:THR:CG2	2:B:128:ASP:OD1	2.55	0.54
2:H:201:MET:HG3	2:H:302:TYR:CE1	2.41	0.54
2:L:63:ALA:HB1	2:L:67:ARG:HD2	1.87	0.54
1:M:6:ILE:HD12	1:M:23:ALA:HB3	1.89	0.54
2:N:444:TYR:HA	2:N:471:GLY:CA	2.31	0.54
1:K:26:SER:OG	1:K:108:LEU:HB3	2.07	0.54
1:O:26:SER:OG	1:O:108:LEU:HB3	2.07	0.54
2:F:298:LEU:HD23	2:F:300:LEU:HD11	1.88	0.54
2:H:104:VAL:HG21	2:H:128:ASP:CB	2.37	0.54
2:P:365:VAL:HG23	2:P:387:VAL:HA	1.89	0.54
2:B:404:ASN:HA	2:B:437:LYS:HD2	1.89	0.54
2:F:357:GLY:CA	2:F:415:ARG:HH22	2.12	0.54
2:H:363:GLY:O	2:H:364:LEU:C	2.46	0.54
2:P:454:GLY:O	2:P:457:TYR:HB2	2.08	0.54
1:I:137:PHE:CD1	2:J:17:VAL:HG21	2.43	0.54
2:N:95:ASN:O	2:N:582:PRO:HG3	2.07	0.54
2:L:91:LEU:O	2:L:567:HIS:HE1	1.89	0.54
1:C:134:ARG:HB2	1:C:139:ILE:O	2.07	0.54
2:N:533:MET:HE3	2:N:588:LEU:HD13	1.89	0.54
1:E:52:LEU:HD12	1:E:52:LEU:O	2.06	0.54
2:H:375:GLN:HG2	2:H:401:TYR:CE1	2.42	0.54
2:N:314:ILE:HD11	2:N:329:ILE:HD11	1.90	0.54
1:A:158:ALA:HA	2:B:62:THR:CG2	2.37	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:N:465:VAL:HG11	2:N:468:MET:HG3	1.89	0.54
2:J:477:ASP:OD2	2:J:504:ALA:HB2	2.08	0.54
2:L:350:GLU:HB3	3:V:209:LEU:HD21	1.89	0.54
2:J:455:LEU:HD22	2:J:483:PHE:HB2	1.88	0.54
2:H:54:HIS:ND1	2:H:55:VAL:N	2.56	0.54
2:P:357:GLY:CA	2:P:415:ARG:HH22	2.12	0.54
2:J:367:GLN:HG3	2:J:390:ILE:HA	1.89	0.54
1:M:99:PHE:CD2	2:N:16:THR:C	2.81	0.54
2:L:547:SER:HB3	2:L:564:HIS:CG	2.43	0.54
2:F:547:SER:HB3	2:F:564:HIS:HB3	1.88	0.54
2:H:428:VAL:HG11	2:H:443:PHE:CE2	2.42	0.54
2:P:227:VAL:HG13	2:P:228:GLY:N	2.22	0.54
2:D:468:MET:HE3	2:D:470:LEU:HD21	1.90	0.54
1:A:145:PRO:HB2	2:D:539:TYR:CE2	2.42	0.54
1:M:159:PHE:O	1:M:160:GLU:CB	2.47	0.54
1:M:160:GLU:HA	1:M:160:GLU:OE2	2.08	0.54
2:N:446:ARG:HG2	2:N:447:GLN:H	1.73	0.54
2:P:367:GLN:HB3	2:P:391:THR:CG2	2.34	0.54
2:H:506:ALA:HB1	2:H:535:MET:HB2	1.88	0.54
2:F:519:TRP:CZ3	2:F:567:HIS:CG	2.95	0.54
2:J:440:ARG:HB3	2:J:467:TRP:CE3	2.43	0.54
1:A:58:TYR:CE1	1:A:62:HIS:CE1	2.96	0.54
2:N:211:ASP:O	2:N:215:ILE:HG13	2.08	0.54
2:P:344:LEU:HD23	2:P:380:LEU:HD21	1.90	0.54
2:P:121:ARG:NH2	5:P:1103:PO4:O4	2.40	0.54
2:P:152:THR:HG22	2:P:177:SER:HB2	1.90	0.54
2:H:153:THR:CG2	2:H:178:GLU:HA	2.38	0.54
1:A:137:PHE:HD1	2:B:17:VAL:HG21	1.72	0.54
2:N:78:LEU:HD12	2:N:79:LYS:H	1.72	0.54
2:H:266:LEU:HD13	2:H:267:VAL:N	2.22	0.54
2:J:357:GLY:CA	2:J:415:ARG:HH22	2.13	0.54
1:A:160:GLU:OE2	1:A:160:GLU:HA	2.08	0.54
2:N:55:VAL:CG2	2:N:75:LEU:HD21	2.37	0.54
2:L:153:THR:HG23	2:L:178:GLU:HA	1.90	0.54
2:H:478:GLU:O	2:H:482:GLU:HG2	2.06	0.54
1:C:93:ILE:HD12	1:C:97:THR:CG2	2.32	0.54
2:D:521:GLN:HG3	2:D:567:HIS:HD2	1.71	0.54
1:E:128:LYS:HB3	1:E:132:GLU:HB2	1.91	0.54
2:L:490:LEU:HD11	2:L:493:LEU:HD13	1.90	0.54
2:L:261:GLU:HG2	2:L:264:MET:HG3	1.90	0.54
2:N:54:HIS:HE1	2:N:56:THR:OG1	1.91	0.54
2:N:357:GLY:CA	2:N:415:ARG:HH22	2.14	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:N:320:LEU:HD21	2:N:323:LEU:HB2	1.90	0.53
2:L:54:HIS:HE1	2:L:56:THR:OG1	1.90	0.53
2:J:227:VAL:CG1	2:J:228:GLY:N	2.71	0.53
2:P:143:LEU:HD23	2:P:159:ILE:HD13	1.90	0.53
2:B:521:GLN:HG3	2:B:567:HIS:CD2	2.43	0.53
2:F:80:LEU:HB2	2:F:122:MET:HE1	1.90	0.53
1:I:48:THR:HG22	1:I:51:ILE:HB	1.91	0.53
1:C:6:ILE:HD12	1:C:23:ALA:HB3	1.90	0.53
2:P:221:SER:O	2:P:223:VAL:HG23	2.08	0.53
1:E:105:ALA:HB3	1:E:114:LEU:CD1	2.32	0.53
2:N:143:LEU:CD2	2:N:159:ILE:HD13	2.38	0.53
2:L:101:THR:OG1	2:L:102:PRO:HD3	2.08	0.53
1:K:134:ARG:HB2	1:K:139:ILE:O	2.09	0.53
2:J:201:MET:HB3	3:U:211:ARG:HH12	1.74	0.53
2:F:65:PRO:HA	2:F:103:TRP:CZ3	2.44	0.53
2:L:314:ILE:HD11	2:L:329:ILE:HD11	1.90	0.53
2:F:85:ARG:NH1	5:F:1101:PO4:O3	2.41	0.53
1:K:113:LEU:O	1:K:117:THR:HG23	2.08	0.53
2:D:270:ARG:HB3	1:G:107:TYR:CD1	2.43	0.53
1:C:48:THR:HG22	1:C:51:ILE:CG1	2.38	0.53
1:I:6:ILE:HG22	1:I:7:VAL:N	2.24	0.53
2:B:501:SER:HA	2:B:524:ARG:HB2	1.91	0.53
1:A:98:LEU:HD21	1:A:120:THR:HG22	1.90	0.53
1:M:91:MET:O	1:M:93:ILE:N	2.40	0.53
1:K:91:MET:O	1:K:93:ILE:N	2.40	0.53
2:N:367:GLN:HG3	2:N:390:ILE:HA	1.90	0.53
2:B:294:GLN:NE2	1:O:107:TYR:OH	2.42	0.53
2:H:98:GLY:O	2:H:122:MET:HE3	2.08	0.53
2:J:547:SER:HB3	2:J:564:HIS:CG	2.44	0.53
2:L:465:VAL:HG11	2:L:468:MET:HG3	1.91	0.53
2:B:298:LEU:HB2	2:B:320:LEU:HD11	1.90	0.53
2:H:43:ARG:HH21	2:H:47:ILE:HD11	1.73	0.53
2:J:399:GLY:O	2:J:434:GLY:HA3	2.08	0.53
2:D:55:VAL:HG23	2:D:75:LEU:HD21	1.91	0.53
2:N:454:GLY:O	2:N:457:TYR:HB2	2.09	0.53
2:F:404:ASN:HA	2:F:437:LYS:HD2	1.89	0.53
2:H:111:LEU:C	2:H:113:GLN:H	2.10	0.53
1:C:91:MET:O	1:C:93:ILE:N	2.41	0.53
2:F:444:TYR:HA	2:F:471:GLY:CA	2.32	0.53
2:B:367:GLN:HB3	2:B:391:THR:CG2	2.34	0.53
2:B:184:LEU:HD12	2:B:207:ILE:HB	1.90	0.53
2:J:468:MET:HE3	2:J:470:LEU:HD21	1.91	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:153:ARG:HG2	1:I:157:TRP:CZ3	2.44	0.53
2:L:121:ARG:NH2	5:L:1103:PO4:O4	2.42	0.53
1:M:87:ASP:OD2	1:M:116:LEU:HD22	2.09	0.53
2:H:364:LEU:HD22	2:H:388:SER:N	2.24	0.53
2:F:367:GLN:HG3	2:F:390:ILE:HA	1.91	0.53
2:L:367:GLN:HG3	2:L:390:ILE:HA	1.89	0.53
2:H:423:PRO:HB3	2:H:425:ASP:OD2	2.09	0.53
1:C:134:ARG:HH11	2:D:40:VAL:HA	1.74	0.53
2:F:397:SER:O	2:F:400:THR:HG22	2.08	0.53
2:B:80:LEU:HB2	2:B:122:MET:HE1	1.91	0.53
2:D:63:ALA:HB1	2:D:67:ARG:HD2	1.91	0.53
1:A:159:PHE:O	1:A:160:GLU:CB	2.52	0.53
2:H:286:PRO:C	2:H:288:LEU:N	2.62	0.53
2:H:362:GLU:C	2:H:364:LEU:N	2.61	0.53
2:P:325:THR:O	2:P:349:ILE:HA	2.09	0.53
2:D:101:THR:CG2	2:D:128:ASP:OD1	2.56	0.53
2:F:311:CYS:HB3	2:F:336:VAL:HG21	1.91	0.53
2:F:153:THR:HG23	2:F:178:GLU:HA	1.91	0.53
1:K:160:GLU:OE2	1:K:160:GLU:HA	2.09	0.53
1:E:11:SER:HA	1:E:45:PRO:HA	1.90	0.53
4:H:4100:OGK:C18	4:H:4100:OGK:HN08	2.21	0.53
1:M:102:ILE:HB	2:N:20:VAL:CG2	2.39	0.53
2:N:519:TRP:CZ3	2:N:567:HIS:CG	2.97	0.53
1:E:26:SER:OG	1:E:108:LEU:HB3	2.09	0.53
2:L:440:ARG:HB3	2:L:467:TRP:CE3	2.43	0.53
2:F:547:SER:HB3	2:F:564:HIS:CG	2.44	0.53
1:A:48:THR:HG22	1:A:51:ILE:CG1	2.39	0.53
2:H:97:GLY:HA3	2:H:123:ILE:CD1	2.38	0.53
2:D:443:PHE:CE2	2:D:445:LEU:HD11	2.44	0.53
2:D:455:LEU:HD22	2:D:483:PHE:HB2	1.90	0.53
1:K:93:ILE:HD12	1:K:97:THR:CG2	2.28	0.53
2:N:93:PRO:HA	2:N:548:ARG:CB	2.33	0.53
2:B:419:ILE:HD11	2:B:446:ARG:HH22	1.74	0.53
2:B:367:GLN:HG3	2:B:390:ILE:HA	1.90	0.53
2:H:65:PRO:HG3	2:H:103:TRP:CD2	2.44	0.53
2:N:59:LEU:O	2:N:62:THR:HB	2.08	0.53
1:O:158:ALA:HA	2:P:62:THR:CG2	2.39	0.53
1:M:134:ARG:CZ	1:M:141:ASN:HD22	2.22	0.53
2:J:227:VAL:HG13	2:J:228:GLY:N	2.23	0.53
2:H:214:THR:HA	2:H:217:ARG:HG2	1.91	0.53
2:N:197:LEU:O	2:N:225:VAL:HA	2.09	0.53
2:L:80:LEU:HB2	2:L:122:MET:CE	2.39	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:93:PRO:HA	2:B:548:ARG:CB	2.32	0.52
1:G:160:GLU:HB3	2:H:31:PRO:HB3	1.90	0.52
2:H:187:LEU:HB3	2:H:215:ILE:HD11	1.90	0.52
2:H:111:LEU:O	2:H:113:GLN:N	2.42	0.52
1:I:58:TYR:CD2	1:I:113:LEU:HD13	2.44	0.52
2:B:275:LEU:HD11	2:B:288:LEU:HD21	1.89	0.52
2:J:422:LEU:HB3	2:J:423:PRO:HD3	1.90	0.52
2:H:184:LEU:HD12	2:H:207:ILE:HB	1.91	0.52
2:L:542:ILE:HD11	2:L:588:LEU:CD1	2.24	0.52
2:H:280:MET:HE1	2:H:285:MET:HA	1.90	0.52
2:H:535:MET:CE	2:H:542:ILE:HG21	2.39	0.52
1:O:58:TYR:CD2	1:O:113:LEU:HD13	2.44	0.52
1:O:6:ILE:HD12	1:O:23:ALA:HB3	1.90	0.52
2:B:399:GLY:O	2:B:434:GLY:HA3	2.09	0.52
2:H:494:GLU:HG2	2:H:519:TRP:HB3	1.91	0.52
2:P:411:VAL:HG22	2:P:444:TYR:HB3	1.91	0.52
1:M:52:LEU:HD12	1:M:52:LEU:O	2.09	0.52
2:N:101:THR:HG22	2:N:128:ASP:HB3	1.90	0.52
2:H:492:LYS:HZ3	2:H:516:ARG:HH11	1.57	0.52
1:I:134:ARG:CZ	1:I:141:ASN:HD22	2.23	0.52
2:L:397:SER:O	2:L:400:THR:HG22	2.08	0.52
2:F:454:GLY:O	2:F:457:TYR:HB2	2.09	0.52
2:N:138:ASP:OD2	2:N:164:ARG:HG3	2.09	0.52
1:I:159:PHE:O	1:I:160:GLU:CB	2.49	0.52
1:M:102:ILE:CD1	2:N:20:VAL:CG2	2.88	0.52
2:P:546:PRO:HD3	2:P:584:THR:O	2.08	0.52
1:M:99:PHE:HB2	2:N:15:ALA:HB3	1.90	0.52
1:C:11:SER:HA	1:C:45:PRO:HA	1.90	0.52
2:D:190:HIS:ND1	2:H:112:ARG:HD2	2.24	0.52
2:F:337:LEU:HD12	2:F:341:CYS:SG	2.49	0.52
2:H:192:THR:HG22	2:H:218:ASN:O	2.10	0.52
2:D:54:HIS:HE1	2:D:56:THR:OG1	1.92	0.52
2:N:233:LEU:O	2:N:236:VAL:HG23	2.08	0.52
2:L:344:LEU:HD23	2:L:380:LEU:HD21	1.91	0.52
2:D:235:LEU:O	2:D:238:PHE:HB3	2.10	0.52
2:J:456:SER:HB2	2:J:482:GLU:HB3	1.91	0.52
2:H:354:ASP:O	2:H:354:ASP:OD2	2.28	0.52
1:C:160:GLU:OE2	1:C:160:GLU:HA	2.09	0.52
2:H:365:VAL:HG21	2:H:387:VAL:CA	2.39	0.52
2:F:289:PHE:HB2	2:F:290:PRO:HD3	1.91	0.52
2:D:519:TRP:CZ3	2:D:567:HIS:CG	2.97	0.52
2:P:55:VAL:CG2	2:P:75:LEU:HD21	2.37	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:402:LEU:HD13	2:H:405:LEU:HG	1.92	0.52
1:G:48:THR:HG22	1:G:51:ILE:CG1	2.39	0.52
2:J:232:ILE:HD12	2:J:252:SER:H	1.74	0.52
1:O:160:GLU:HA	1:O:160:GLU:OE2	2.10	0.52
2:L:446:ARG:HG2	2:L:447:GLN:H	1.74	0.52
2:D:547:SER:HB3	2:D:564:HIS:CG	2.43	0.52
1:C:26:SER:OG	1:C:108:LEU:HB3	2.10	0.52
2:N:547:SER:HB3	2:N:564:HIS:CG	2.44	0.52
1:O:137:PHE:HD1	2:P:17:VAL:HG21	1.72	0.52
2:P:275:LEU:HD11	2:P:288:LEU:HD21	1.91	0.52
2:P:501:SER:HA	2:P:524:ARG:HB2	1.92	0.52
2:P:446:ARG:HG2	2:P:447:GLN:H	1.74	0.52
2:N:472:TYR:OH	3:W:201:LEU:HB2	2.10	0.52
2:D:446:ARG:HG2	2:D:447:GLN:H	1.74	0.52
2:N:325:THR:O	2:N:349:ILE:HA	2.09	0.52
2:H:170:LEU:HD23	2:H:170:LEU:C	2.30	0.52
2:L:545:ILE:HG22	2:L:546:PRO:CD	2.40	0.52
1:I:158:ALA:HA	2:J:62:THR:HG21	1.90	0.52
2:L:235:LEU:O	2:L:238:PHE:HB3	2.10	0.52
2:F:521:GLN:HG3	2:F:567:HIS:CD2	2.45	0.52
2:L:521:GLN:HG3	2:L:567:HIS:CD2	2.44	0.52
2:D:275:LEU:HD11	2:D:288:LEU:HD21	1.92	0.52
2:N:130:ASP:OD1	2:N:134:LYS:HE3	2.09	0.52
1:K:153:ARG:NH2	2:L:539:TYR:CE1	2.78	0.52
2:D:375:GLN:HG2	2:D:401:TYR:CE1	2.45	0.52
2:N:456:SER:HB2	2:N:482:GLU:HB3	1.92	0.52
2:B:454:GLY:O	2:B:457:TYR:HB2	2.10	0.52
2:F:533:MET:HE3	2:F:588:LEU:HD13	1.91	0.52
2:F:411:VAL:HG22	2:F:444:TYR:HB3	1.92	0.52
2:D:367:GLN:HG3	2:D:390:ILE:HA	1.92	0.52
2:P:419:ILE:CD1	2:P:446:ARG:HH22	2.23	0.52
2:D:85:ARG:NH2	4:D:2100:OGK:O07	2.35	0.52
2:J:521:GLN:HG3	2:J:567:HIS:CD2	2.45	0.52
2:H:225:VAL:CG2	2:H:245:LEU:HD11	2.40	0.52
2:D:311:CYS:HB3	2:D:336:VAL:HG21	1.91	0.52
2:J:275:LEU:HD11	2:J:288:LEU:HD21	1.90	0.52
2:J:465:VAL:HG11	2:J:468:MET:HG3	1.91	0.52
2:B:54:HIS:HE1	2:B:56:THR:OG1	1.92	0.52
2:J:305:LEU:HD23	2:J:305:LEU:O	2.10	0.52
2:H:306:GLU:O	2:H:310:HIS:HD2	1.92	0.52
2:J:93:PRO:HA	2:J:548:ARG:CB	2.33	0.52
4:B:1100:OGK:C18	4:B:1100:OGK:HN08	2.19	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:J:85:ARG:NH2	4:J:5100:OGK:O07	2.33	0.52
2:N:298:LEU:HB2	2:N:320:LEU:HD11	1.92	0.52
2:P:407:ASP:OD1	2:P:440:ARG:HD2	2.10	0.52
2:H:328:VAL:HG12	2:H:359:GLU:HG2	1.92	0.52
2:F:422:LEU:HB3	2:F:423:PRO:HD3	1.92	0.52
2:L:404:ASN:HA	2:L:437:LYS:HD2	1.92	0.52
2:B:337:LEU:HD12	2:B:341:CYS:SG	2.50	0.52
2:J:80:LEU:HB2	2:J:122:MET:CE	2.39	0.52
1:C:137:PHE:CD1	2:D:17:VAL:HG21	2.45	0.52
1:E:93:ILE:HD12	1:E:97:THR:CG2	2.29	0.51
1:G:151:VAL:CG1	2:H:39:LEU:HD21	2.40	0.51
2:P:367:GLN:HG3	2:P:390:ILE:HA	1.92	0.51
2:D:519:TRP:CZ3	2:D:567:HIS:ND1	2.79	0.51
1:M:137:PHE:HD1	2:N:17:VAL:HG21	1.74	0.51
2:N:101:THR:CG2	2:N:128:ASP:OD1	2.58	0.51
2:D:191:ASN:ND2	2:D:194:LEU:H	2.07	0.51
2:N:310:HIS:O	2:N:314:ILE:CG1	2.58	0.51
2:H:302:TYR:N	2:H:302:TYR:CD2	2.74	0.51
2:P:80:LEU:HB2	2:P:122:MET:CE	2.40	0.51
2:H:532:LEU:HD11	2:H:568:ILE:HD13	1.92	0.51
2:P:456:SER:HB2	2:P:482:GLU:HB3	1.92	0.51
2:F:235:LEU:O	2:F:238:PHE:HB3	2.10	0.51
2:H:411:VAL:CG2	4:H:4100:OGK:H16A	2.40	0.51
2:B:367:GLN:O	2:B:371:ILE:HG22	2.09	0.51
1:M:112:ASN:C	1:M:114:LEU:N	2.62	0.51
2:N:456:SER:CB	2:N:482:GLU:HB3	2.40	0.51
2:L:477:ASP:OD2	2:L:504:ALA:HB2	2.10	0.51
2:P:233:LEU:O	2:P:236:VAL:HG23	2.11	0.51
2:P:251:GLY:O	2:P:278:SER:HB2	2.10	0.51
2:B:446:ARG:HG2	2:B:447:GLN:H	1.75	0.51
2:P:101:THR:OG1	2:P:102:PRO:HD3	2.10	0.51
2:P:398:ILE:HG23	2:P:402:LEU:HD11	1.92	0.51
2:F:233:LEU:O	2:F:236:VAL:HG23	2.09	0.51
2:H:418:ARG:HG2	2:H:446:ARG:NH1	2.25	0.51
2:H:455:LEU:HD21	2:H:473:VAL:CG2	2.40	0.51
1:C:158:ALA:HA	2:D:62:THR:HG21	1.91	0.51
2:D:298:LEU:HB2	2:D:320:LEU:HD11	1.93	0.51
1:A:11:SER:HA	1:A:45:PRO:HA	1.92	0.51
1:M:87:ASP:HB3	1:M:116:LEU:HD21	1.92	0.51
2:H:406:CYS:HA	2:H:438:LEU:HA	1.92	0.51
2:J:235:LEU:O	2:J:238:PHE:HB3	2.10	0.51
2:F:63:ALA:HB1	2:F:67:ARG:HD2	1.91	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:455:LEU:HD22	2:B:483:PHE:HB2	1.91	0.51
2:D:93:PRO:HA	2:D:548:ARG:CB	2.32	0.51
2:H:65:PRO:HA	2:H:103:TRP:CZ3	2.44	0.51
2:B:143:LEU:CD2	2:B:159:ILE:HD13	2.40	0.51
1:E:48:THR:HG22	1:E:51:ILE:CG1	2.40	0.51
1:O:48:THR:CG2	1:O:51:ILE:HG12	2.41	0.51
2:N:153:THR:HG23	2:N:178:GLU:HA	1.93	0.51
2:H:424:LEU:O	2:H:426:ASN:N	2.43	0.51
2:N:337:LEU:HD12	2:N:341:CYS:SG	2.51	0.51
2:N:542:ILE:CD1	2:N:588:LEU:HD12	2.28	0.51
2:H:382:TYR:CD2	2:H:407:ASP:HB3	2.46	0.51
1:A:52:LEU:HD12	1:A:52:LEU:O	2.10	0.51
2:F:546:PRO:HD3	2:F:584:THR:O	2.08	0.51
1:A:105:ALA:HB2	1:A:113:LEU:HD23	1.92	0.51
1:M:48:THR:HG22	1:M:51:ILE:CG1	2.40	0.51
2:J:311:CYS:HB3	2:J:336:VAL:HG21	1.91	0.51
1:I:48:THR:HG22	1:I:51:ILE:CG1	2.40	0.51
2:F:255:GLU:HG3	2:F:255:GLU:O	2.11	0.51
2:H:278:SER:C	2:H:280:MET:H	2.13	0.51
2:H:496:ARG:HD2	4:H:4100:OGK:H23A	1.92	0.51
1:A:134:ARG:HH11	2:B:40:VAL:HA	1.75	0.51
2:P:63:ALA:HB1	2:P:67:ARG:HD2	1.91	0.51
1:A:153:ARG:NH2	2:B:539:TYR:CE1	2.79	0.51
2:J:501:SER:HA	2:J:524:ARG:HB2	1.92	0.51
2:D:422:LEU:HB3	2:D:423:PRO:HD3	1.93	0.51
2:P:247:GLU:HA	2:P:274:ARG:O	2.11	0.51
2:H:306:GLU:OE2	2:H:306:GLU:HA	2.09	0.51
2:L:519:TRP:HH2	2:L:567:HIS:CE1	2.29	0.51
2:P:191:ASN:ND2	2:P:194:LEU:H	2.06	0.51
2:F:298:LEU:HB2	2:F:320:LEU:HD11	1.92	0.51
1:K:137:PHE:HD1	2:L:17:VAL:HG21	1.76	0.51
1:K:48:THR:HG22	1:K:51:ILE:CG1	2.41	0.51
2:P:232:ILE:HD12	2:P:252:SER:H	1.76	0.51
2:L:78:LEU:HD12	2:L:79:LYS:H	1.76	0.51
1:O:91:MET:CE	1:O:117:THR:HG22	2.41	0.51
2:F:367:GLN:O	2:F:371:ILE:HG22	2.10	0.51
2:P:78:LEU:HD12	2:P:79:LYS:H	1.74	0.51
2:D:310:HIS:O	2:D:314:ILE:HG12	2.11	0.51
1:O:91:MET:O	1:O:93:ILE:N	2.44	0.51
2:H:414:ASP:HB3	2:H:446:ARG:HH21	1.76	0.51
1:O:11:SER:HA	1:O:45:PRO:HA	1.92	0.51
2:F:272:LEU:HD21	2:F:275:LEU:HB3	1.92	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:272:LEU:HD21	2:B:275:LEU:HB3	1.93	0.51
2:N:422:LEU:HB3	2:N:423:PRO:HD3	1.93	0.51
2:F:501:SER:HA	2:F:524:ARG:HB2	1.93	0.51
2:F:305:LEU:HD23	2:F:305:LEU:O	2.11	0.51
2:H:365:VAL:HG21	2:H:387:VAL:CG1	2.39	0.50
2:H:390:ILE:HD11	2:H:412:LEU:CD2	2.41	0.50
2:H:409:ARG:HD2	4:H:4100:OGK:H13A	1.92	0.50
2:B:519:TRP:HH2	2:B:567:HIS:CE1	2.29	0.50
2:B:546:PRO:HD3	2:B:584:THR:O	2.11	0.50
2:B:320:LEU:HD21	2:B:323:LEU:HB2	1.93	0.50
2:D:272:LEU:HD21	2:D:275:LEU:HB3	1.92	0.50
1:M:153:ARG:HG2	1:M:157:TRP:CZ3	2.46	0.50
2:B:153:THR:HG23	2:B:178:GLU:HA	1.92	0.50
2:B:305:LEU:O	2:B:305:LEU:HD23	2.11	0.50
2:B:255:GLU:HG3	2:B:255:GLU:O	2.11	0.50
2:F:542:ILE:CD1	2:F:588:LEU:HD12	2.30	0.50
2:F:143:LEU:CD2	2:F:159:ILE:HD13	2.41	0.50
1:C:134:ARG:NH1	2:D:40:VAL:HA	2.26	0.50
2:B:482:GLU:O	2:B:485:ARG:HG2	2.11	0.50
2:B:357:GLY:CA	2:B:415:ARG:HH22	2.12	0.50
2:P:289:PHE:HB2	2:P:290:PRO:HD3	1.92	0.50
2:H:370:LEU:HB2	2:H:394:SER:HB3	1.93	0.50
2:F:398:ILE:O	2:F:402:LEU:HD12	2.12	0.50
2:H:291:PHE:N	2:H:291:PHE:CD2	2.79	0.50
2:P:65:PRO:HA	2:P:103:TRP:CZ3	2.46	0.50
2:L:482:GLU:O	2:L:485:ARG:HG2	2.11	0.50
2:F:110:ASN:HA	2:L:190:HIS:HE1	1.77	0.50
2:P:153:THR:HG23	2:P:178:GLU:HA	1.92	0.50
1:C:91:MET:O	1:C:93:ILE:HG12	2.12	0.50
2:H:87:ALA:HB2	2:H:92:ILE:HB	1.93	0.50
1:K:128:LYS:HB3	1:K:132:GLU:HB2	1.93	0.50
2:H:230:PHE:CB	2:H:235:LEU:HD21	2.41	0.50
1:I:158:ALA:HA	2:J:62:THR:CG2	2.41	0.50
2:P:159:ILE:HD12	2:P:166:ILE:HD11	1.93	0.50
2:F:95:ASN:O	2:F:582:PRO:HG3	2.12	0.50
2:D:456:SER:HB2	2:D:482:GLU:HB3	1.94	0.50
2:H:432:LEU:HD11	2:H:458:ILE:HD13	1.93	0.50
2:N:523:TYR:CE2	2:N:568:ILE:HD11	2.47	0.50
2:J:519:TRP:HH2	2:J:567:HIS:CE1	2.30	0.50
2:B:440:ARG:HB3	2:B:467:TRP:CE3	2.45	0.50
2:B:80:LEU:HB2	2:B:122:MET:CE	2.41	0.50
2:B:130:ASP:OD1	2:B:134:LYS:HE3	2.12	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:233:LEU:O	2:L:236:VAL:HG23	2.11	0.50
2:H:145:LEU:O	2:H:171:MET:HA	2.10	0.50
2:J:419:ILE:HD11	2:J:446:ARG:HH22	1.75	0.50
1:K:105:ALA:HB2	1:K:113:LEU:HD23	1.93	0.50
2:D:20:VAL:O	2:D:24:VAL:HG23	2.11	0.50
2:P:422:LEU:HB3	2:P:423:PRO:HD3	1.93	0.50
2:H:464:ASN:O	2:H:466:ARG:HD2	2.11	0.50
2:P:289:PHE:HD1	2:P:316:LYS:HD2	1.73	0.50
2:J:289:PHE:HB2	2:J:290:PRO:HD3	1.93	0.50
1:A:128:LYS:HB3	1:A:132:GLU:HB2	1.94	0.50
2:B:159:ILE:HD12	2:B:166:ILE:HD11	1.94	0.50
1:E:48:THR:HG22	1:E:51:ILE:CB	2.41	0.50
2:F:477:ASP:OD2	2:F:504:ALA:HB2	2.12	0.50
2:H:392:ASN:ND2	2:H:422:LEU:HD13	2.26	0.50
1:K:26:SER:HB3	1:K:29:ILE:HG13	1.93	0.50
1:I:128:LYS:HB3	1:I:132:GLU:HB2	1.94	0.50
2:F:465:VAL:HG11	2:F:468:MET:HG3	1.93	0.50
2:J:398:ILE:HG23	2:J:402:LEU:HD11	1.94	0.50
2:H:289:PHE:N	2:H:290:PRO:CD	2.74	0.50
2:H:282:PRO:HB3	2:H:309:ASP:OD2	2.11	0.50
2:H:496:ARG:HB2	4:H:4100:OGK:H01A	1.94	0.50
2:H:367:GLN:N	2:H:367:GLN:CD	2.65	0.50
2:L:320:LEU:HD21	2:L:323:LEU:HB2	1.93	0.50
1:M:130:PRO:O	1:M:134:ARG:HG2	2.12	0.50
2:P:482:GLU:O	2:P:485:ARG:HG2	2.12	0.50
2:N:227:VAL:CG1	2:N:228:GLY:N	2.74	0.50
2:F:197:LEU:O	2:F:225:VAL:HA	2.12	0.50
2:B:232:ILE:HD12	2:B:252:SER:H	1.76	0.50
2:F:232:ILE:HD12	2:F:252:SER:H	1.76	0.50
1:G:113:LEU:O	1:G:117:THR:HG23	2.11	0.50
1:K:91:MET:O	1:K:93:ILE:HG12	2.12	0.49
2:P:419:ILE:HD11	2:P:446:ARG:HH22	1.75	0.49
2:J:164:ARG:HH21	2:N:136:ARG:HH21	1.60	0.49
2:D:407:ASP:OD1	2:D:440:ARG:HD2	2.12	0.49
2:H:191:ASN:ND2	2:H:194:LEU:H	2.08	0.49
1:O:113:LEU:HG	1:O:113:LEU:O	2.12	0.49
1:G:11:SER:HA	1:G:45:PRO:HA	1.94	0.49
2:H:516:ARG:HA	2:H:572:TYR:HD1	1.75	0.49
2:P:465:VAL:HG11	2:P:468:MET:HG3	1.94	0.49
2:B:310:HIS:O	2:B:314:ILE:CG1	2.58	0.49
2:J:80:LEU:HB2	2:J:122:MET:HE1	1.94	0.49
2:P:80:LEU:HB2	2:P:122:MET:HE1	1.94	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:310:HIS:O	2:D:314:ILE:CG1	2.60	0.49
2:J:54:HIS:HE1	2:J:56:THR:OG1	1.95	0.49
2:H:289:PHE:N	2:H:290:PRO:HD2	2.27	0.49
1:M:102:ILE:CB	2:N:20:VAL:CG2	2.90	0.49
2:N:565:PRO:HG2	3:W:201:LEU:HD21	1.94	0.49
2:H:191:ASN:ND2	2:H:193:SER:H	2.10	0.49
1:O:134:ARG:CZ	1:O:141:ASN:HD22	2.25	0.49
2:D:398:ILE:HG23	2:D:402:LEU:HD11	1.94	0.49
2:L:310:HIS:O	2:L:314:ILE:HG12	2.12	0.49
2:H:97:GLY:CA	2:H:123:ILE:HD11	2.42	0.49
2:P:477:ASP:OD2	2:P:504:ALA:HB2	2.12	0.49
1:E:91:MET:O	1:E:93:ILE:N	2.45	0.49
2:H:387:VAL:HG11	2:H:390:ILE:CD1	2.39	0.49
1:K:132:GLU:O	1:K:136:THR:HG23	2.12	0.49
2:J:519:TRP:CZ3	2:J:567:HIS:CG	3.01	0.49
2:H:357:GLY:C	2:H:359:GLU:N	2.66	0.49
1:C:134:ARG:CZ	1:C:141:ASN:HD22	2.24	0.49
2:P:138:ASP:OD2	2:P:164:ARG:HG3	2.12	0.49
2:N:357:GLY:HA2	2:N:415:ARG:NH2	2.14	0.49
2:L:519:TRP:CZ3	2:L:567:HIS:CG	3.00	0.49
2:B:350:GLU:HB3	3:Q:209:LEU:CD2	2.40	0.49
1:C:58:TYR:CE1	1:C:62:HIS:CE1	3.00	0.49
2:P:468:MET:CE	2:P:470:LEU:HD11	2.42	0.49
1:A:134:ARG:NH1	2:B:40:VAL:HA	2.27	0.49
2:P:159:ILE:HG13	2:P:160:VAL:N	2.26	0.49
2:D:65:PRO:HA	2:D:103:TRP:CZ3	2.48	0.49
2:N:116:SER:CB	2:N:142:THR:HG23	2.36	0.49
1:A:91:MET:O	1:A:93:ILE:HG12	2.13	0.49
1:I:91:MET:O	1:I:93:ILE:N	2.42	0.49
2:H:487:CYS:HB3	2:H:490:LEU:HB2	1.95	0.49
2:B:519:TRP:CZ3	2:B:567:HIS:CG	3.00	0.49
2:H:492:LYS:HZ3	2:H:516:ARG:NH1	2.09	0.49
2:B:419:ILE:CD1	2:B:446:ARG:HH22	2.26	0.49
2:J:411:VAL:HG22	2:J:444:TYR:HB3	1.93	0.49
1:G:10:SER:OG	1:G:11:SER:N	2.46	0.49
2:N:159:ILE:HG13	2:N:160:VAL:N	2.27	0.49
2:H:428:VAL:CG1	2:H:443:PHE:CZ	2.96	0.49
1:M:48:THR:HG22	1:M:51:ILE:CB	2.42	0.49
2:J:159:ILE:HD12	2:J:166:ILE:HD11	1.94	0.49
1:K:134:ARG:HH11	2:L:40:VAL:HA	1.76	0.49
2:L:80:LEU:HB2	2:L:122:MET:HE1	1.95	0.49
2:N:375:GLN:HG2	2:N:401:TYR:CE1	2.48	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:211:ASP:O	2:D:215:ILE:HG13	2.12	0.49
2:J:404:ASN:HA	2:J:437:LYS:HD2	1.94	0.49
2:J:443:PHE:CE2	2:J:445:LEU:HD11	2.48	0.49
2:P:248:PHE:CD2	2:P:248:PHE:C	2.86	0.49
1:A:26:SER:OG	1:A:108:LEU:HB3	2.13	0.49
1:O:128:LYS:HB3	1:O:132:GLU:HB2	1.94	0.49
2:N:289:PHE:HB2	2:N:290:PRO:HD3	1.94	0.49
2:D:519:TRP:HH2	2:D:567:HIS:CE1	2.31	0.49
1:C:128:LYS:HB3	1:C:132:GLU:HB2	1.95	0.49
2:F:295:ILE:HG21	2:F:298:LEU:CD1	2.41	0.49
2:P:320:LEU:HD21	2:P:323:LEU:HB2	1.93	0.49
2:F:407:ASP:OD1	2:F:440:ARG:HD2	2.12	0.49
2:H:83:LYS:O	2:H:121:ARG:NH1	2.35	0.49
2:N:235:LEU:O	2:N:238:PHE:HB3	2.12	0.49
2:L:223:VAL:O	2:L:245:LEU:HD12	2.13	0.49
2:H:332:ARG:O	2:H:336:VAL:HG23	2.13	0.49
2:H:450:LEU:HD11	2:H:454:GLY:CA	2.39	0.49
2:H:450:LEU:CG	2:H:454:GLY:HA3	2.43	0.49
2:B:546:PRO:O	2:B:547:SER:HB2	2.13	0.49
2:J:490:LEU:HD11	2:J:493:LEU:HD13	1.94	0.49
1:O:10:SER:OG	1:O:11:SER:N	2.46	0.49
2:L:227:VAL:HG13	2:L:228:GLY:H	1.76	0.49
2:N:400:THR:O	2:N:403:LYS:HE2	2.12	0.49
1:O:153:ARG:HG2	1:O:157:TRP:CZ3	2.48	0.49
2:J:482:GLU:O	2:J:485:ARG:HG2	2.13	0.49
2:H:431:LEU:HD23	2:H:432:LEU:N	2.27	0.49
2:L:255:GLU:HG3	2:L:255:GLU:O	2.13	0.49
2:N:170:LEU:HD23	2:N:170:LEU:C	2.33	0.49
2:P:255:GLU:O	2:P:255:GLU:HG3	2.13	0.49
2:H:153:THR:HG23	2:H:178:GLU:HA	1.95	0.49
2:F:409:ARG:HD3	4:F:3100:OGK:H13A	1.93	0.49
1:I:134:ARG:HH11	2:J:40:VAL:HA	1.77	0.49
2:L:275:LEU:HD11	2:L:288:LEU:HD21	1.93	0.49
2:D:482:GLU:O	2:D:485:ARG:HG2	2.12	0.49
2:D:247:GLU:HA	2:D:274:ARG:O	2.12	0.49
2:P:130:ASP:OD1	2:P:134:LYS:HE3	2.12	0.49
2:N:469:LEU:HA	2:N:494:GLU:O	2.11	0.49
2:H:385:VAL:HG12	2:H:387:VAL:HG23	1.95	0.49
2:H:259:MET:N	2:H:260:PRO:HD3	2.28	0.49
2:H:57:MET:HB2	2:H:80:LEU:HD23	1.95	0.49
2:N:311:CYS:CB	2:N:336:VAL:HG21	2.43	0.49
2:L:298:LEU:HB2	2:L:320:LEU:HD11	1.94	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:490:LEU:HD11	2:F:493:LEU:HD13	1.95	0.49
2:J:153:THR:HG23	2:J:178:GLU:HA	1.94	0.49
1:K:159:PHE:O	1:K:160:GLU:CB	2.50	0.48
2:D:57:MET:CE	2:D:62:THR:HG22	2.43	0.48
2:N:546:PRO:O	2:N:547:SER:HB2	2.12	0.48
2:H:43:ARG:NH2	2:H:47:ILE:HD11	2.28	0.48
1:K:134:ARG:NH1	2:L:40:VAL:HA	2.28	0.48
2:J:397:SER:O	2:J:400:THR:HG22	2.13	0.48
2:L:538:PRO:O	2:L:539:TYR:HB2	2.13	0.48
2:N:96:TRP:O	2:N:578:ARG:NH2	2.41	0.48
2:D:153:THR:HG23	2:D:178:GLU:HA	1.95	0.48
2:N:152:THR:HG22	2:N:177:SER:HB2	1.95	0.48
2:F:538:PRO:O	2:F:539:TYR:HB2	2.13	0.48
2:B:321:GLU:HA	2:B:344:LEU:HA	1.94	0.48
2:L:546:PRO:HD3	2:L:584:THR:O	2.11	0.48
2:N:57:MET:CE	2:N:62:THR:HG22	2.43	0.48
1:K:58:TYR:CD2	1:K:113:LEU:HD13	2.47	0.48
2:J:400:THR:O	2:J:403:LYS:HE2	2.13	0.48
2:N:482:GLU:O	2:N:485:ARG:HG2	2.13	0.48
2:L:501:SER:HA	2:L:524:ARG:HB2	1.95	0.48
2:H:526:SER:O	2:H:527:MET:C	2.52	0.48
2:H:326:ARG:O	2:H:329:ILE:HG22	2.12	0.48
2:H:85:ARG:NH1	5:H:1101:PO4:O2	2.45	0.48
2:B:282:PRO:HA	2:B:285:MET:CE	2.44	0.48
2:L:419:ILE:CD1	2:L:446:ARG:HH22	2.27	0.48
2:P:519:TRP:CZ3	2:P:567:HIS:CG	3.01	0.48
2:F:519:TRP:CZ3	2:F:567:HIS:ND1	2.81	0.48
1:A:48:THR:CG2	1:A:51:ILE:HG12	2.43	0.48
2:B:176:PHE:CZ	2:B:204:PHE:CZ	3.01	0.48
2:F:247:GLU:HA	2:F:274:ARG:O	2.13	0.48
2:J:255:GLU:HG3	2:J:255:GLU:O	2.12	0.48
2:D:255:GLU:O	2:D:255:GLU:HG3	2.13	0.48
2:H:454:GLY:O	2:H:457:TYR:HB2	2.12	0.48
1:M:128:LYS:HB2	1:M:133:ILE:CD1	2.43	0.48
2:N:419:ILE:CD1	2:N:446:ARG:HH22	2.26	0.48
1:E:134:ARG:HH11	2:F:40:VAL:HA	1.78	0.48
2:L:519:TRP:CZ3	2:L:567:HIS:ND1	2.82	0.48
2:J:546:PRO:HD3	2:J:584:THR:O	2.14	0.48
1:O:48:THR:HG22	1:O:51:ILE:CB	2.43	0.48
2:D:397:SER:O	2:D:400:THR:HG22	2.13	0.48
2:P:456:SER:CB	2:P:482:GLU:HB3	2.43	0.48
2:F:231:GLU:HG2	2:F:254:ASN:HD22	1.78	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:152:THR:HG22	2:F:177:SER:HB2	1.95	0.48
2:L:578:ARG:HG3	2:L:578:ARG:H	1.44	0.48
2:B:247:GLU:HA	2:B:274:ARG:O	2.14	0.48
2:D:399:GLY:O	2:D:434:GLY:HA3	2.13	0.48
2:F:46:LYS:HE3	2:F:46:LYS:HB2	1.61	0.48
2:H:486:GLY:O	2:H:487:CYS:C	2.51	0.48
2:L:85:ARG:NH1	5:L:1101:PO4:O3	2.46	0.48
2:J:519:TRP:CZ3	2:J:567:HIS:ND1	2.82	0.48
2:N:159:ILE:HD12	2:N:166:ILE:HD11	1.95	0.48
1:C:153:ARG:NH2	2:D:539:TYR:CE1	2.81	0.48
2:H:476:SER:C	2:H:478:GLU:H	2.16	0.48
2:B:539:TYR:OH	1:C:146:GLU:HA	2.13	0.48
2:D:454:GLY:O	2:D:457:TYR:HB2	2.14	0.48
1:I:111:LYS:O	1:I:115:ASP:HB2	2.12	0.48
2:L:454:GLY:O	2:L:457:TYR:HB2	2.13	0.48
2:N:501:SER:HA	2:N:524:ARG:HB2	1.95	0.48
2:D:296:ARG:NH2	1:G:37:CYS:SG	2.87	0.48
2:D:305:LEU:HD23	2:D:305:LEU:O	2.13	0.48
2:J:454:GLY:O	2:J:457:TYR:HB2	2.14	0.48
2:H:364:LEU:HD21	2:H:386:TYR:O	2.13	0.48
2:J:546:PRO:O	2:J:547:SER:HB2	2.13	0.48
2:H:391:THR:C	2:H:393:GLU:N	2.67	0.48
2:J:295:ILE:HG21	2:J:298:LEU:CD1	2.43	0.48
1:G:137:PHE:HB3	2:H:17:VAL:CG2	2.43	0.48
2:N:80:LEU:HB2	2:N:122:MET:HE2	1.95	0.48
1:G:48:THR:CG2	1:G:51:ILE:HG12	2.44	0.48
2:L:398:ILE:O	2:L:402:LEU:HD12	2.12	0.48
2:H:316:LYS:O	2:H:318:PRO:HD3	2.14	0.48
2:F:446:ARG:HG2	2:F:447:GLN:N	2.28	0.48
2:D:546:PRO:HD3	2:D:584:THR:O	2.13	0.48
2:F:101:THR:HG22	2:F:128:ASP:HB3	1.95	0.48
2:B:477:ASP:N	2:B:477:ASP:OD1	2.46	0.48
2:J:456:SER:CB	2:J:482:GLU:HB3	2.43	0.48
2:H:431:LEU:CD2	2:H:432:LEU:HD23	2.43	0.48
1:E:153:ARG:HG2	1:E:157:TRP:CZ3	2.49	0.48
2:F:127:LEU:HD21	2:F:131:ARG:NH2	2.28	0.48
2:B:170:LEU:HD23	2:B:170:LEU:C	2.34	0.48
1:I:26:SER:HB3	1:I:29:ILE:HG13	1.96	0.48
2:N:419:ILE:HD11	2:N:446:ARG:HH22	1.79	0.48
2:L:289:PHE:HD1	2:L:316:LYS:HD2	1.73	0.48
2:H:542:ILE:HD11	2:H:588:LEU:HD12	1.96	0.48
2:H:57:MET:HE2	2:H:62:THR:HG22	1.95	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:M:11:SER:HA	1:M:45:PRO:HA	1.96	0.48
2:N:545:ILE:HG22	2:N:546:PRO:CD	2.44	0.48
1:A:105:ALA:HB3	1:A:114:LEU:HD13	1.94	0.48
2:D:465:VAL:HG11	2:D:468:MET:HG3	1.96	0.48
2:L:477:ASP:OD1	2:L:477:ASP:N	2.47	0.48
2:L:456:SER:HB2	2:L:482:GLU:HB3	1.94	0.48
2:J:121:ARG:NH2	5:J:1103:PO4:O4	2.47	0.48
2:P:176:PHE:CZ	2:P:204:PHE:CZ	3.02	0.48
2:H:21:ILE:HD12	2:H:24:VAL:HB	1.95	0.48
2:N:282:PRO:HA	2:N:285:MET:CE	2.43	0.48
1:E:134:ARG:CZ	1:E:141:ASN:HD22	2.26	0.48
2:H:492:LYS:HZ1	2:H:516:ARG:NH1	2.10	0.48
2:N:392:ASN:HD21	2:N:424:LEU:HA	1.77	0.48
2:L:251:GLY:O	2:L:278:SER:HB2	2.13	0.48
2:D:176:PHE:CZ	2:D:204:PHE:CZ	3.01	0.48
2:H:365:VAL:HG21	2:H:387:VAL:HA	1.95	0.48
2:L:419:ILE:HD11	2:L:446:ARG:HH22	1.79	0.48
2:P:521:GLN:HG3	2:P:567:HIS:CD2	2.48	0.48
2:H:170:LEU:HB2	2:H:198:ASN:HB3	1.95	0.48
2:H:477:ASP:OD2	2:H:501:SER:OG	2.23	0.48
1:A:48:THR:HG22	1:A:51:ILE:CB	2.43	0.48
1:G:6:ILE:HG22	1:G:7:VAL:N	2.28	0.48
2:N:387:VAL:O	2:N:413:LEU:HD22	2.14	0.48
2:F:523:TYR:CE2	2:F:568:ILE:HD11	2.49	0.48
1:K:11:SER:HA	1:K:45:PRO:HA	1.95	0.48
2:F:96:TRP:O	2:F:578:ARG:NH2	2.42	0.48
2:D:321:GLU:HA	2:D:344:LEU:HA	1.96	0.48
2:L:127:LEU:HD21	2:L:131:ARG:NH2	2.29	0.48
2:P:170:LEU:HD23	2:P:170:LEU:C	2.34	0.48
1:E:45:PRO:CB	2:L:291:PHE:HA	2.44	0.47
2:H:412:LEU:HD12	2:H:445:LEU:CA	2.44	0.47
2:L:367:GLN:O	2:L:371:ILE:HG22	2.14	0.47
2:D:367:GLN:O	2:D:371:ILE:HG22	2.14	0.47
2:B:519:TRP:CZ3	2:B:567:HIS:ND1	2.82	0.47
2:J:545:ILE:HG22	2:J:546:PRO:CD	2.43	0.47
2:L:298:LEU:HD23	2:L:300:LEU:HD11	1.95	0.47
1:E:48:THR:HG23	1:E:51:ILE:H	1.79	0.47
1:O:134:ARG:HH11	2:P:40:VAL:HA	1.78	0.47
2:N:191:ASN:ND2	2:N:194:LEU:H	2.12	0.47
2:N:443:PHE:CE2	2:N:445:LEU:HD11	2.48	0.47
2:H:364:LEU:HD22	2:H:365:VAL:HG22	1.96	0.47
2:J:164:ARG:HD2	2:N:112:ARG:CZ	2.45	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:N:519:TRP:CZ3	2:N:567:HIS:ND1	2.82	0.47
2:N:289:PHE:HD1	2:N:316:LYS:HD2	1.76	0.47
2:J:298:LEU:HB2	2:J:320:LEU:HD11	1.96	0.47
2:D:295:ILE:HG21	2:D:298:LEU:CD1	2.43	0.47
2:B:227:VAL:HG13	2:B:228:GLY:H	1.78	0.47
2:N:351:ARG:O	2:N:351:ARG:HG3	2.14	0.47
1:I:48:THR:CG2	1:I:51:ILE:HG12	2.43	0.47
2:L:456:SER:CB	2:L:482:GLU:HB3	2.44	0.47
2:H:247:GLU:HA	2:H:274:ARG:O	2.14	0.47
2:H:114:LEU:O	2:H:136:ARG:NH1	2.48	0.47
2:H:64:THR:O	2:H:65:PRO:C	2.51	0.47
2:H:204:PHE:O	2:H:206:LYS:N	2.47	0.47
2:N:490:LEU:HD11	2:N:493:LEU:HD13	1.96	0.47
1:M:48:THR:CG2	1:M:51:ILE:HG12	2.45	0.47
1:G:120:THR:O	1:G:124:MET:HG3	2.14	0.47
1:C:48:THR:CG2	1:C:51:ILE:HG12	2.44	0.47
2:D:477:ASP:OD2	2:D:504:ALA:HB2	2.12	0.47
2:J:63:ALA:HB1	2:J:67:ARG:HD2	1.96	0.47
2:N:171:MET:O	2:N:174:SER:HB2	2.14	0.47
2:J:247:GLU:HA	2:J:274:ARG:O	2.14	0.47
2:D:232:ILE:HD12	2:D:252:SER:H	1.79	0.47
2:N:176:PHE:CZ	2:N:204:PHE:CZ	3.02	0.47
2:D:412:LEU:HD12	2:D:412:LEU:C	2.35	0.47
2:H:280:MET:HG2	2:H:280:MET:O	2.14	0.47
2:J:446:ARG:HG2	2:J:447:GLN:N	2.29	0.47
2:F:297:LYS:HE3	2:F:297:LYS:HB2	1.65	0.47
2:L:545:ILE:HG22	2:L:546:PRO:N	2.30	0.47
2:N:297:LYS:HB2	2:N:297:LYS:HE3	1.62	0.47
2:D:565:PRO:HG2	3:R:201:LEU:HD21	1.97	0.47
2:P:440:ARG:HB3	2:P:467:TRP:CE3	2.45	0.47
1:I:11:SER:HA	1:I:45:PRO:HA	1.95	0.47
2:N:247:GLU:HA	2:N:274:ARG:O	2.14	0.47
2:D:501:SER:HA	2:D:524:ARG:HB2	1.96	0.47
2:N:248:PHE:CD2	2:N:248:PHE:C	2.87	0.47
2:B:248:PHE:CD2	2:B:248:PHE:C	2.87	0.47
1:E:105:ALA:HB2	1:E:113:LEU:HD23	1.96	0.47
2:H:89:PHE:CE2	4:H:4100:OGK:H04A	2.50	0.47
2:D:289:PHE:HD1	2:D:316:LYS:HD2	1.73	0.47
1:E:132:GLU:O	1:E:136:THR:HG23	2.15	0.47
2:B:465:VAL:HG11	2:B:468:MET:HG3	1.95	0.47
1:K:134:ARG:CZ	1:K:141:ASN:HD22	2.28	0.47
2:J:398:ILE:C	2:J:400:THR:H	2.17	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:138:ASP:OD2	2:F:164:ARG:HG3	2.14	0.47
2:B:422:LEU:HB3	2:B:423:PRO:HD3	1.96	0.47
2:H:315:GLN:HG3	2:H:340:TYR:CE1	2.50	0.47
1:I:52:LEU:HD12	1:I:52:LEU:O	2.14	0.47
2:H:262:LYS:HD2	2:H:263:TYR:CE1	2.49	0.47
2:D:80:LEU:HB2	2:D:122:MET:HE1	1.96	0.47
2:P:477:ASP:OD1	2:P:477:ASP:N	2.48	0.47
2:P:543:GLU:OE2	2:P:578:ARG:HD3	2.13	0.47
2:H:269:PRO:O	2:H:270:ARG:C	2.52	0.47
2:F:375:GLN:HG2	2:F:401:TYR:CE1	2.49	0.47
2:H:53:GLU:O	2:H:75:LEU:HD22	2.15	0.47
2:H:153:THR:HG23	2:H:177:SER:O	2.15	0.47
2:H:52:ARG:NH1	2:H:72:PHE:HZ	2.11	0.47
2:J:285:MET:N	2:J:286:PRO:CD	2.78	0.47
2:P:546:PRO:O	2:P:547:SER:HB2	2.15	0.47
1:C:26:SER:HB3	1:C:29:ILE:HG13	1.96	0.47
2:D:320:LEU:HD21	2:D:323:LEU:HB2	1.97	0.47
1:C:52:LEU:HD12	1:C:52:LEU:O	2.14	0.47
2:J:57:MET:CE	2:J:62:THR:HG22	2.43	0.47
2:P:398:ILE:O	2:P:402:LEU:HD12	2.14	0.47
2:H:532:LEU:O	2:H:534:GLN:N	2.48	0.47
2:L:543:GLU:OE2	2:L:578:ARG:HD3	2.15	0.47
2:L:339:GLN:HA	2:L:342:LYS:HE2	1.96	0.47
2:L:247:GLU:HA	2:L:274:ARG:O	2.13	0.47
2:D:95:ASN:O	2:D:582:PRO:HG3	2.15	0.47
2:P:444:TYR:HA	2:P:471:GLY:CA	2.37	0.47
1:E:26:SER:HB3	1:E:29:ILE:HG13	1.97	0.47
1:I:35:ASP:CB	2:N:243:ALA:HB1	2.42	0.47
2:H:367:GLN:CA	2:H:367:GLN:OE1	2.62	0.47
2:P:298:LEU:HB2	2:P:320:LEU:HD11	1.96	0.47
2:B:468:MET:CE	2:B:470:LEU:HD11	2.45	0.47
2:P:321:GLU:HA	2:P:344:LEU:HA	1.97	0.47
1:G:105:ALA:HB2	1:G:113:LEU:HD23	1.97	0.47
2:J:472:TYR:CD2	2:J:497:GLY:O	2.68	0.47
2:F:113:GLN:HE22	2:L:192:THR:HG21	1.80	0.47
2:H:352:GLY:O	2:H:353:ALA:HB2	2.15	0.47
2:J:223:VAL:O	2:J:245:LEU:HD12	2.14	0.47
1:M:132:GLU:O	1:M:136:THR:HG23	2.14	0.47
2:N:20:VAL:O	2:N:24:VAL:HG23	2.15	0.47
2:J:289:PHE:HD1	2:J:316:LYS:HD2	1.77	0.47
2:H:468:MET:HG3	2:H:490:LEU:HD21	1.96	0.47
2:J:325:THR:O	2:J:349:ILE:HA	2.15	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:289:PHE:HB2	2:D:290:PRO:HD3	1.97	0.47
2:H:187:LEU:HA	2:H:187:LEU:HD23	1.72	0.47
2:D:190:HIS:HE1	2:H:110:ASN:HA	1.80	0.47
2:F:400:THR:O	2:F:403:LYS:HE2	2.15	0.47
1:C:48:THR:HG22	1:C:51:ILE:CB	2.44	0.47
1:G:58:TYR:CE1	1:G:62:HIS:CE1	3.03	0.47
2:H:168:THR:HB	2:H:196:VAL:HG13	1.96	0.47
2:H:58:ALA:HA	2:H:81:LYS:HD2	1.97	0.47
2:L:297:LYS:HB2	2:L:297:LYS:HE3	1.64	0.47
2:F:440:ARG:HB3	2:F:467:TRP:CE3	2.48	0.47
1:I:134:ARG:NH1	2:J:40:VAL:HA	2.30	0.47
1:C:114:LEU:C	1:C:116:LEU:H	2.17	0.47
2:J:310:HIS:O	2:J:314:ILE:CG1	2.63	0.47
1:I:91:MET:O	1:I:93:ILE:HG12	2.14	0.46
2:H:284:GLU:O	2:H:287:ILE:HG23	2.15	0.46
1:G:134:ARG:HH11	2:H:40:VAL:HA	1.80	0.46
1:M:137:PHE:HD1	2:N:17:VAL:CG2	2.27	0.46
2:F:320:LEU:HD21	2:F:323:LEU:HB2	1.96	0.46
2:L:546:PRO:O	2:L:547:SER:HB2	2.15	0.46
2:J:191:ASN:ND2	2:J:194:LEU:H	2.11	0.46
2:H:197:LEU:O	2:H:225:VAL:HA	2.15	0.46
2:H:230:PHE:HD1	2:H:235:LEU:HD21	1.78	0.46
2:D:477:ASP:N	2:D:477:ASP:OD1	2.47	0.46
2:L:523:TYR:CE2	2:L:568:ILE:HD11	2.50	0.46
2:H:78:LEU:CD1	2:H:79:LYS:H	2.11	0.46
2:B:357:GLY:HA2	2:B:415:ARG:NH2	2.12	0.46
2:H:31:PRO:O	2:H:35:ASP:HB2	2.16	0.46
1:K:108:LEU:HD12	1:K:110:ILE:CD1	2.40	0.46
2:H:92:ILE:HG23	2:H:93:PRO:HD2	1.97	0.46
1:M:128:LYS:HB3	1:M:132:GLU:HB2	1.96	0.46
2:P:519:TRP:CZ3	2:P:567:HIS:ND1	2.83	0.46
2:P:519:TRP:HH2	2:P:567:HIS:CE1	2.32	0.46
2:F:159:ILE:HG13	2:F:160:VAL:N	2.29	0.46
2:B:590:GLU:O	2:B:591:PRO:C	2.54	0.46
2:P:85:ARG:NH1	5:P:1101:PO4:O3	2.48	0.46
2:P:311:CYS:CB	2:P:336:VAL:HG21	2.46	0.46
1:E:6:ILE:HG22	1:E:7:VAL:N	2.30	0.46
2:F:396:GLU:HG2	2:F:430:SER:OG	2.15	0.46
2:J:396:GLU:HG2	2:J:430:SER:OG	2.16	0.46
1:C:114:LEU:C	1:C:116:LEU:N	2.68	0.46
2:P:373:LEU:HD12	2:P:377:CYS:SG	2.54	0.46
2:L:65:PRO:HA	2:L:103:TRP:CZ3	2.50	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:91:MET:O	1:G:93:ILE:N	2.47	0.46
2:H:469:LEU:HA	2:H:494:GLU:O	2.14	0.46
2:P:285:MET:N	2:P:286:PRO:CD	2.79	0.46
2:B:545:ILE:HG22	2:B:546:PRO:CD	2.45	0.46
2:N:490:LEU:O	2:N:514:SER:HB2	2.14	0.46
1:M:58:TYR:CE1	1:M:62:HIS:CE1	3.04	0.46
1:E:137:PHE:HD1	2:F:17:VAL:HG21	1.79	0.46
2:B:456:SER:HB2	2:B:482:GLU:HB3	1.98	0.46
2:N:227:VAL:HG13	2:N:228:GLY:N	2.30	0.46
2:F:266:LEU:HD13	2:F:266:LEU:HA	1.81	0.46
2:N:338:ALA:O	2:N:376:GLY:HA3	2.15	0.46
2:H:148:CYS:O	2:H:174:SER:HA	2.16	0.46
1:C:159:PHE:O	1:C:160:GLU:CB	2.52	0.46
2:H:450:LEU:HD21	2:H:455:LEU:N	2.30	0.46
1:E:42:VAL:HG22	1:E:42:VAL:O	2.15	0.46
2:B:85:ARG:NH1	5:B:1101:PO4:O3	2.49	0.46
2:L:446:ARG:HG2	2:L:447:GLN:N	2.31	0.46
1:G:26:SER:C	1:G:28:THR:H	2.19	0.46
1:M:99:PHE:CB	2:N:15:ALA:HB1	2.45	0.46
2:L:143:LEU:CD2	2:L:159:ILE:HD13	2.43	0.46
2:N:398:ILE:O	2:N:402:LEU:HD12	2.16	0.46
1:G:48:THR:HG23	1:G:51:ILE:H	1.80	0.46
1:G:48:THR:HG22	1:G:51:ILE:CB	2.45	0.46
2:N:468:MET:CE	2:N:470:LEU:HD11	2.46	0.46
2:L:310:HIS:O	2:L:314:ILE:CG1	2.64	0.46
2:P:235:LEU:O	2:P:238:PHE:HB3	2.14	0.46
2:J:310:HIS:O	2:J:314:ILE:HG12	2.16	0.46
2:L:422:LEU:HB3	2:L:423:PRO:HD3	1.97	0.46
2:J:373:LEU:HD12	2:J:377:CYS:SG	2.55	0.46
2:L:232:ILE:HD12	2:L:252:SER:H	1.80	0.46
2:J:419:ILE:CD1	2:J:446:ARG:HH22	2.28	0.46
2:D:282:PRO:HA	2:D:285:MET:CE	2.45	0.46
2:D:285:MET:N	2:D:286:PRO:CD	2.79	0.46
1:I:106:ASN:ND2	2:J:23:GLN:NE2	2.63	0.46
4:F:3100:OGK:N08	4:F:3100:OGK:C18	2.77	0.46
2:B:270:ARG:O	1:O:107:TYR:HE1	1.98	0.46
2:H:298:LEU:HD13	2:H:300:LEU:HD11	1.96	0.46
2:F:191:ASN:ND2	2:F:194:LEU:H	2.13	0.46
1:G:125:ILE:HD11	2:H:44:TRP:CH2	2.51	0.46
1:C:153:ARG:HG2	1:C:157:TRP:CZ3	2.50	0.46
2:H:29:THR:HG23	2:H:30:ASP:N	2.31	0.46
2:L:46:LYS:HE3	2:L:46:LYS:HB2	1.60	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:160:GLU:HG3	2:H:52:ARG:NH2	2.18	0.46
2:L:367:GLN:CB	2:L:391:THR:HG22	2.39	0.46
2:P:57:MET:CE	2:P:62:THR:HG22	2.46	0.46
2:F:456:SER:CB	2:F:482:GLU:HB3	2.46	0.46
2:D:299:ASP:OD2	2:D:301:LEU:HB2	2.16	0.46
2:F:344:LEU:HD23	2:F:380:LEU:HD21	1.98	0.46
2:J:46:LYS:HE3	2:J:46:LYS:HB2	1.55	0.46
2:L:198:ASN:C	2:L:198:ASN:OD1	2.53	0.46
2:J:367:GLN:O	2:J:371:ILE:HG22	2.15	0.46
2:B:367:GLN:CB	2:B:391:THR:HG22	2.41	0.46
1:K:52:LEU:O	1:K:52:LEU:HD12	2.16	0.46
1:O:134:ARG:NH1	2:P:40:VAL:HA	2.31	0.46
2:B:539:TYR:CE2	1:C:145:PRO:HB2	2.51	0.46
2:J:294:GLN:NE2	1:M:107:TYR:OH	2.49	0.46
1:M:46:ASN:HB2	1:M:107:TYR:CZ	2.51	0.46
2:P:282:PRO:HA	2:P:285:MET:CE	2.43	0.46
2:D:419:ILE:HD11	2:D:446:ARG:HH22	1.81	0.46
2:B:289:PHE:N	2:B:290:PRO:CD	2.79	0.46
2:P:545:ILE:HG22	2:P:546:PRO:CD	2.45	0.46
2:B:253:LEU:CD1	2:B:280:MET:HB2	2.42	0.46
2:F:101:THR:N	2:F:102:PRO:CD	2.79	0.46
2:B:191:ASN:ND2	2:B:194:LEU:H	2.11	0.46
2:L:101:THR:CG2	2:L:128:ASP:OD1	2.63	0.46
2:B:398:ILE:O	2:B:402:LEU:HD12	2.16	0.46
2:H:291:PHE:N	2:H:291:PHE:HD2	2.11	0.46
1:C:137:PHE:HD1	2:D:17:VAL:HG21	1.79	0.46
2:F:121:ARG:NH2	5:F:1103:PO4:O4	2.48	0.46
1:O:159:PHE:O	1:O:160:GLU:CB	2.51	0.46
2:H:253:LEU:HD12	2:H:280:MET:HB2	1.97	0.46
2:H:285:MET:HE1	2:H:309:ASP:HB3	1.98	0.46
2:H:390:ILE:CD1	2:H:410:LEU:HD11	2.45	0.46
2:L:325:THR:O	2:L:349:ILE:HA	2.16	0.46
2:P:446:ARG:HG2	2:P:447:GLN:N	2.31	0.46
2:H:501:SER:HB2	2:H:503:ARG:NH2	2.31	0.46
2:J:538:PRO:O	2:J:539:TYR:HB2	2.16	0.46
2:F:392:ASN:HD21	2:F:424:LEU:HA	1.80	0.46
2:L:248:PHE:CD2	2:L:248:PHE:C	2.89	0.46
2:P:542:ILE:CD1	2:P:588:LEU:HD12	2.32	0.45
1:G:26:SER:HB3	1:G:29:ILE:HG13	1.97	0.45
2:H:80:LEU:HD11	2:H:103:TRP:CD2	2.51	0.45
2:N:545:ILE:HG22	2:N:546:PRO:HD2	1.98	0.45
2:N:546:PRO:HB2	2:N:547:SER:H	1.67	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:102:ILE:CG2	1:I:118:CYS:SG	3.03	0.45
2:F:468:MET:CE	2:F:470:LEU:HD11	2.46	0.45
2:H:129:LEU:HD23	2:H:155:GLY:HA3	1.98	0.45
2:H:233:LEU:HD21	2:H:262:LYS:HG2	1.97	0.45
2:D:412:LEU:O	2:D:412:LEU:HD12	2.15	0.45
2:J:20:VAL:O	2:J:24:VAL:HG23	2.16	0.45
2:N:339:GLN:HA	2:N:342:LYS:HE2	1.98	0.45
2:F:170:LEU:C	2:F:170:LEU:HD23	2.36	0.45
2:D:116:SER:CB	2:D:142:THR:HG23	2.40	0.45
2:H:365:VAL:CG2	2:H:388:SER:H	2.29	0.45
2:N:496:ARG:NH2	3:W:201:LEU:HD22	2.31	0.45
2:B:289:PHE:HD1	2:B:316:LYS:HD2	1.75	0.45
2:F:398:ILE:HG23	2:F:402:LEU:HD11	1.97	0.45
1:I:58:TYR:CE1	1:I:62:HIS:CE1	3.04	0.45
2:B:351:ARG:O	2:B:351:ARG:HG3	2.16	0.45
2:B:327:ASN:HD22	2:B:327:ASN:N	2.13	0.45
2:B:436:LYS:HB3	2:B:436:LYS:HE2	1.72	0.45
1:G:130:PRO:HG3	2:H:39:LEU:HB2	1.98	0.45
2:P:409:ARG:HD3	4:P:8100:OGK:H13A	1.98	0.45
1:G:52:LEU:O	1:G:52:LEU:HD12	2.16	0.45
2:F:387:VAL:O	2:F:413:LEU:HD22	2.17	0.45
2:D:468:MET:CE	2:D:470:LEU:HD11	2.46	0.45
2:J:176:PHE:CZ	2:J:204:PHE:CZ	3.04	0.45
5:N:1102:PO4:O3	3:W:206:ARG:NH2	2.45	0.45
2:F:436:LYS:HE2	2:F:436:LYS:HB3	1.72	0.45
1:E:159:PHE:O	1:E:160:GLU:CB	2.51	0.45
1:G:93:ILE:HD12	1:G:97:THR:CG2	2.36	0.45
4:L:6100:OGK:N08	4:L:6100:OGK:C18	2.79	0.45
2:B:101:THR:HG22	2:B:128:ASP:CB	2.46	0.45
2:F:490:LEU:HD23	2:F:490:LEU:HA	1.74	0.45
2:P:387:VAL:O	2:P:413:LEU:HD22	2.17	0.45
2:D:96:TRP:O	2:D:578:ARG:NH2	2.42	0.45
2:F:351:ARG:O	2:F:351:ARG:HG3	2.15	0.45
2:L:305:LEU:HD23	2:L:305:LEU:O	2.16	0.45
1:M:91:MET:SD	1:M:117:THR:HG22	2.55	0.45
2:H:418:ARG:HB3	2:H:421:ASP:HB2	1.99	0.45
2:F:367:GLN:CB	2:F:391:THR:HG22	2.41	0.45
2:H:487:CYS:HB2	2:H:512:LEU:CD2	2.46	0.45
2:F:472:TYR:CD2	2:F:497:GLY:O	2.70	0.45
2:J:546:PRO:O	2:J:547:SER:CB	2.65	0.45
1:M:137:PHE:CD1	2:N:17:VAL:CG2	2.99	0.45
2:H:243:ALA:C	2:H:245:LEU:H	2.20	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:503:ARG:HG2	2:H:504:ALA:N	2.31	0.45
1:M:6:ILE:HG22	1:M:7:VAL:N	2.31	0.45
2:D:456:SER:CB	2:D:482:GLU:HB3	2.47	0.45
2:P:523:TYR:CE2	2:P:568:ILE:HD11	2.52	0.45
2:B:65:PRO:HA	2:B:103:TRP:CZ3	2.52	0.45
1:E:58:TYR:CE1	1:E:62:HIS:CE1	3.04	0.45
2:B:338:ALA:O	2:B:376:GLY:HA3	2.17	0.45
1:E:19:GLU:N	1:E:19:GLU:OE1	2.47	0.45
2:N:255:GLU:O	2:N:255:GLU:HG3	2.16	0.45
1:M:91:MET:O	1:M:93:ILE:HG12	2.16	0.45
1:O:91:MET:O	1:O:93:ILE:HG12	2.17	0.45
1:E:10:SER:OG	1:E:11:SER:N	2.48	0.45
2:F:519:TRP:HH2	2:F:567:HIS:CE1	2.34	0.45
1:E:134:ARG:NH1	2:F:40:VAL:HA	2.32	0.45
2:L:159:ILE:HD12	2:L:166:ILE:HD11	1.98	0.45
1:E:48:THR:CG2	1:E:51:ILE:HG12	2.46	0.45
2:L:57:MET:CE	2:L:62:THR:HG22	2.47	0.45
1:O:48:THR:HG23	1:O:51:ILE:H	1.82	0.45
2:P:285:MET:N	2:P:286:PRO:HD2	2.32	0.45
1:I:125:ILE:HG23	1:I:133:ILE:CD1	2.39	0.45
2:N:446:ARG:HG2	2:N:447:GLN:N	2.32	0.45
2:D:446:ARG:HG2	2:D:447:GLN:N	2.32	0.45
2:H:351:ARG:HG3	2:H:359:GLU:OE2	2.16	0.45
2:F:456:SER:HB2	2:F:482:GLU:HB3	1.98	0.45
2:N:440:ARG:HB3	2:N:467:TRP:CE3	2.51	0.45
1:C:42:VAL:HA	1:C:43:PRO:HD3	1.79	0.45
2:D:127:LEU:HB2	2:N:126:ASP:HB3	1.99	0.45
2:J:127:LEU:HD21	2:J:131:ARG:NH2	2.32	0.45
2:L:563:GLU:O	2:L:563:GLU:HG3	2.17	0.45
2:D:327:ASN:HD22	2:D:327:ASN:N	2.15	0.45
2:D:357:GLY:CA	2:D:415:ARG:HH22	2.14	0.45
2:H:387:VAL:HG12	2:H:390:ILE:HD13	1.95	0.45
1:M:102:ILE:CB	2:N:20:VAL:HG21	2.47	0.45
2:L:282:PRO:HA	2:L:285:MET:CE	2.47	0.45
2:B:545:ILE:HB	2:B:567:HIS:HB2	1.98	0.45
2:B:289:PHE:HB2	2:B:290:PRO:HD3	1.97	0.45
2:L:419:ILE:O	2:L:420:THR:O	2.34	0.45
1:I:102:ILE:HG12	1:I:117:THR:CB	2.47	0.45
3:X:201:LEU:HD23	3:X:201:LEU:HA	1.80	0.45
1:K:6:ILE:HG22	1:K:7:VAL:N	2.32	0.45
2:D:400:THR:O	2:D:403:LYS:HE2	2.17	0.45
2:B:351:ARG:HD3	2:B:413:LEU:HD11	1.98	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:46:ASN:HB2	1:I:107:TYR:CZ	2.52	0.45
2:J:187:LEU:HD23	2:J:187:LEU:HA	1.82	0.45
2:D:248:PHE:CD2	2:D:248:PHE:C	2.90	0.45
2:H:91:LEU:O	2:H:567:HIS:HE1	2.00	0.45
2:D:289:PHE:N	2:D:290:PRO:CD	2.80	0.45
1:G:42:VAL:O	1:G:42:VAL:HG22	2.17	0.45
1:I:102:ILE:HG12	1:I:117:THR:HB	1.99	0.45
1:O:137:PHE:HD1	2:P:17:VAL:CG2	2.29	0.45
2:N:407:ASP:OD1	2:N:440:ARG:HD2	2.17	0.45
2:D:523:TYR:CE2	2:D:568:ILE:HD11	2.52	0.45
2:J:198:ASN:OD1	2:J:198:ASN:C	2.55	0.45
2:L:544:LEU:HD11	2:L:588:LEU:HD11	1.99	0.45
2:H:400:THR:C	2:H:403:LYS:HE2	2.37	0.45
1:G:42:VAL:HA	1:G:43:PRO:HD3	1.79	0.45
1:M:99:PHE:CZ	2:N:17:VAL:HG22	2.52	0.45
2:F:546:PRO:O	2:F:547:SER:HB2	2.17	0.45
2:L:55:VAL:CG2	2:L:75:LEU:HD21	2.44	0.45
2:D:159:ILE:HD12	2:D:166:ILE:HD11	1.99	0.45
1:C:6:ILE:HG22	1:C:7:VAL:N	2.32	0.45
2:H:96:TRP:HE3	2:H:97:GLY:N	2.15	0.45
1:O:6:ILE:HG22	1:O:7:VAL:N	2.32	0.45
2:H:53:GLU:O	2:H:75:LEU:CD2	2.65	0.45
2:D:351:ARG:O	2:D:351:ARG:HG3	2.16	0.45
2:N:429:ARG:O	2:N:433:ILE:HD12	2.17	0.45
2:F:89:PHE:CD1	3:S:206:ARG:HA	2.52	0.45
2:L:412:LEU:HD12	2:L:412:LEU:C	2.37	0.45
2:H:275:LEU:C	2:H:275:LEU:HD12	2.37	0.45
2:H:467:TRP:HZ3	2:H:494:GLU:HG3	1.82	0.44
2:N:367:GLN:CB	2:N:391:THR:HG22	2.37	0.44
2:J:285:MET:N	2:J:286:PRO:HD2	2.32	0.44
2:D:546:PRO:O	2:D:547:SER:HB2	2.16	0.44
2:J:490:LEU:HD23	2:J:490:LEU:HA	1.79	0.44
2:D:490:LEU:HD11	2:D:493:LEU:HD13	1.98	0.44
2:B:311:CYS:CB	2:B:336:VAL:HG21	2.47	0.44
2:N:398:ILE:HG23	2:N:402:LEU:HD11	1.99	0.44
1:K:48:THR:CG2	1:K:51:ILE:HG12	2.48	0.44
2:P:120:ARG:NH2	5:P:1103:PO4:O4	2.48	0.44
2:D:578:ARG:HG3	2:D:578:ARG:H	1.46	0.44
2:F:251:GLY:O	2:F:278:SER:HB2	2.17	0.44
2:F:20:VAL:O	2:F:24:VAL:HG23	2.18	0.44
1:K:42:VAL:HA	1:K:43:PRO:HD3	1.81	0.44
2:P:339:GLN:HA	2:P:342:LYS:HE2	1.99	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:412:LEU:HD12	2:B:412:LEU:C	2.37	0.44
2:F:544:LEU:HD11	2:F:588:LEU:HD11	2.00	0.44
2:J:419:ILE:O	2:J:420:THR:C	2.55	0.44
2:B:419:ILE:O	2:B:420:THR:O	2.35	0.44
3:W:201:LEU:HA	3:W:202:PRO:HD3	1.83	0.44
2:P:367:GLN:CB	2:P:391:THR:HG22	2.42	0.44
2:L:468:MET:CE	2:L:470:LEU:HD11	2.48	0.44
2:B:59:LEU:O	2:B:62:THR:HB	2.17	0.44
2:J:321:GLU:HA	2:J:344:LEU:HA	1.99	0.44
2:D:398:ILE:C	2:D:400:THR:H	2.20	0.44
2:H:97:GLY:HA3	2:H:123:ILE:HD11	1.98	0.44
1:E:153:ARG:NH2	2:F:539:TYR:CE1	2.85	0.44
2:L:443:PHE:CE2	2:L:445:LEU:HD11	2.52	0.44
2:L:373:LEU:HD12	2:L:377:CYS:SG	2.57	0.44
2:P:375:GLN:HG2	2:P:401:TYR:CE1	2.52	0.44
2:B:392:ASN:HD21	2:B:424:LEU:HA	1.82	0.44
2:N:46:LYS:HE3	2:N:46:LYS:HB2	1.54	0.44
1:O:91:MET:HE3	1:O:117:THR:HG22	1.98	0.44
2:H:409:ARG:HA	2:H:442:ALA:O	2.18	0.44
2:H:410:LEU:HD13	2:H:411:VAL:O	2.18	0.44
2:H:87:ALA:HB2	2:H:92:ILE:CG1	2.47	0.44
2:N:519:TRP:HH2	2:N:567:HIS:CE1	2.35	0.44
2:D:419:ILE:CD1	2:D:446:ARG:HH22	2.29	0.44
2:P:545:ILE:HG22	2:P:546:PRO:HD2	1.99	0.44
2:D:521:GLN:HG3	2:D:567:HIS:CD2	2.50	0.44
1:O:26:SER:HB3	1:O:29:ILE:HG13	1.98	0.44
1:O:58:TYR:CE1	1:O:62:HIS:CE1	3.05	0.44
2:F:545:ILE:HG22	2:F:546:PRO:CD	2.47	0.44
2:N:80:LEU:HB2	2:N:122:MET:HE1	1.97	0.44
2:P:96:TRP:O	2:P:578:ARG:NH2	2.44	0.44
2:P:168:THR:CB	2:P:196:VAL:HG13	2.26	0.44
1:E:107:TYR:CD1	2:L:270:ARG:HB3	2.52	0.44
2:F:419:ILE:CD1	2:F:446:ARG:HH22	2.31	0.44
2:B:546:PRO:HB2	2:B:547:SER:H	1.68	0.44
4:D:2100:OGK:N08	4:D:2100:OGK:H18A	2.26	0.44
2:D:545:ILE:HB	2:D:567:HIS:HB2	1.99	0.44
1:A:95:GLN:NE2	1:A:124:MET:HE3	2.32	0.44
2:P:101:THR:HG22	2:P:128:ASP:CB	2.47	0.44
2:L:191:ASN:ND2	2:L:194:LEU:H	2.12	0.44
2:H:423:PRO:HA	2:H:449:GLY:O	2.18	0.44
2:J:380:LEU:HD23	2:J:380:LEU:HA	1.79	0.44
2:P:396:GLU:HG2	2:P:430:SER:OG	2.16	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:400:THR:O	2:L:403:LYS:HE2	2.18	0.44
2:D:351:ARG:HD3	2:D:413:LEU:HD11	1.98	0.44
2:L:338:ALA:O	2:L:376:GLY:HA3	2.17	0.44
2:D:533:MET:CE	2:D:588:LEU:HD13	2.47	0.44
2:J:419:ILE:O	2:J:420:THR:O	2.36	0.44
2:H:208:SER:HA	2:H:209:PRO:HD2	1.82	0.44
2:H:101:THR:HB	2:H:102:PRO:HD3	1.98	0.44
1:O:137:PHE:CD1	2:P:17:VAL:CG2	3.01	0.44
2:J:539:TYR:OH	1:K:146:GLU:HA	2.18	0.44
2:H:97:GLY:CA	2:H:123:ILE:CD1	2.96	0.44
1:I:148:GLU:HG2	1:I:152:ARG:HH12	1.83	0.44
2:J:338:ALA:O	2:J:376:GLY:HA3	2.17	0.44
2:J:304:LEU:HA	2:J:304:LEU:HD12	1.71	0.44
2:P:419:ILE:O	2:P:420:THR:C	2.56	0.44
2:D:419:ILE:O	2:D:420:THR:C	2.56	0.44
2:H:535:MET:HG2	2:H:535:MET:O	2.17	0.44
4:J:5100:OGK:C18	4:J:5100:OGK:N08	2.79	0.44
2:J:85:ARG:NH1	5:J:1101:PO4:O3	2.50	0.44
2:D:297:LYS:HE3	2:D:297:LYS:HB2	1.68	0.44
2:P:314:ILE:HD11	2:P:329:ILE:HD11	1.99	0.44
2:P:538:PRO:O	2:P:539:TYR:HB2	2.18	0.44
2:P:65:PRO:HB3	2:P:103:TRP:CE3	2.52	0.44
2:D:344:LEU:HD23	2:D:380:LEU:HD21	1.98	0.44
2:J:299:ASP:OD2	2:J:301:LEU:HB2	2.18	0.44
2:J:392:ASN:HD21	2:J:424:LEU:HA	1.82	0.44
2:N:231:GLU:HG2	2:N:254:ASN:HD22	1.82	0.44
2:H:511:LYS:HB2	2:H:511:LYS:HE2	1.66	0.44
2:P:419:ILE:O	2:P:420:THR:O	2.36	0.44
2:N:419:ILE:O	2:N:420:THR:C	2.56	0.44
1:M:96:ALA:CB	2:N:14:VAL:CG1	2.95	0.44
4:P:8100:OGK:C18	4:P:8100:OGK:N08	2.75	0.44
1:M:99:PHE:HB2	2:N:15:ALA:HB1	2.00	0.44
2:L:545:ILE:HG22	2:L:546:PRO:HD2	1.98	0.44
2:H:225:VAL:HG21	2:H:238:PHE:CZ	2.53	0.44
1:A:137:PHE:HD1	2:B:17:VAL:CG2	2.31	0.44
2:H:423:PRO:CB	2:H:425:ASP:OD2	2.65	0.44
2:F:55:VAL:CG2	2:F:75:LEU:HD21	2.48	0.44
1:K:48:THR:HG22	1:K:51:ILE:CB	2.47	0.44
2:N:382:TYR:CD2	2:N:407:ASP:HB3	2.53	0.44
2:D:75:LEU:HA	2:D:75:LEU:HD23	1.81	0.44
1:I:153:ARG:NH2	2:J:539:TYR:CE1	2.86	0.44
2:L:321:GLU:HA	2:L:344:LEU:HA	1.98	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:275:LEU:C	2:H:275:LEU:CD1	2.86	0.44
2:N:83:LYS:O	2:N:121:ARG:NH1	2.46	0.44
2:J:263:TYR:O	2:J:266:LEU:HD23	2.18	0.44
2:N:85:ARG:NH1	5:N:1101:PO4:O3	2.51	0.44
2:F:419:ILE:O	2:F:420:THR:C	2.56	0.44
2:B:545:ILE:HG22	2:B:546:PRO:HD2	2.00	0.44
1:E:129:THR:HB	1:E:130:PRO:HD2	2.00	0.44
2:D:319:ASN:HD21	1:G:43:PRO:HG3	1.82	0.44
2:N:101:THR:N	2:N:102:PRO:CD	2.81	0.44
2:F:230:PHE:O	2:F:252:SER:HB3	2.18	0.44
2:D:127:LEU:HD21	2:D:131:ARG:NH2	2.33	0.44
1:G:148:GLU:HG3	1:G:152:ARG:NH1	2.33	0.44
2:D:266:LEU:HD13	2:D:266:LEU:HA	1.80	0.44
2:J:78:LEU:HD12	2:J:79:LYS:H	1.83	0.44
2:L:276:GLY:HA3	2:L:299:ASP:HB3	2.00	0.44
2:J:543:GLU:OE2	2:J:578:ARG:HD3	2.16	0.44
2:D:46:LYS:HB2	2:D:46:LYS:HE3	1.55	0.44
2:H:289:PHE:CE1	2:H:316:LYS:HD3	2.53	0.44
2:H:519:TRP:CH2	2:H:567:HIS:ND1	2.86	0.44
1:C:58:TYR:CD2	1:C:113:LEU:HD13	2.52	0.44
1:O:52:LEU:O	1:O:52:LEU:HD12	2.17	0.44
2:L:75:LEU:HD23	2:L:75:LEU:HA	1.83	0.44
2:H:101:THR:N	2:H:102:PRO:CD	2.81	0.44
2:J:468:MET:CE	2:J:470:LEU:HD11	2.48	0.44
1:I:34:GLU:O	2:N:271:LYS:HE3	2.18	0.44
2:L:176:PHE:CZ	2:L:204:PHE:CZ	3.06	0.44
2:H:189:GLN:C	2:H:190:HIS:ND1	2.71	0.44
2:B:419:ILE:O	2:B:420:THR:C	2.56	0.43
2:B:446:ARG:HG2	2:B:447:GLN:N	2.33	0.43
2:H:390:ILE:HD11	2:H:412:LEU:HD21	2.00	0.43
2:H:411:VAL:HG13	2:H:444:TYR:CB	2.40	0.43
1:I:108:LEU:HD12	1:I:110:ILE:CD1	2.43	0.43
2:D:165:LYS:NZ	2:H:139:ASP:OD1	2.43	0.43
2:J:545:ILE:HG22	2:J:546:PRO:HD2	1.99	0.43
2:P:125:SER:O	2:P:129:LEU:HD22	2.18	0.43
2:N:184:LEU:HA	2:N:184:LEU:HD23	1.86	0.43
1:C:48:THR:HG23	1:C:51:ILE:H	1.83	0.43
2:N:398:ILE:C	2:N:400:THR:H	2.20	0.43
2:D:538:PRO:O	2:D:539:TYR:HB2	2.18	0.43
2:B:18:ASP:OD2	2:B:43:ARG:HD2	2.17	0.43
2:N:251:GLY:O	2:N:278:SER:HB2	2.18	0.43
2:N:305:LEU:HD23	2:N:305:LEU:O	2.18	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:N:436:LYS:HE2	2:N:436:LYS:HB3	1.75	0.43
2:F:590:GLU:O	2:F:591:PRO:C	2.57	0.43
2:F:419:ILE:O	2:F:420:THR:O	2.36	0.43
2:H:512:LEU:HA	2:H:513:PRO:HD3	1.71	0.43
2:J:320:LEU:HD21	2:J:323:LEU:HB2	1.99	0.43
1:I:10:SER:OG	1:I:11:SER:N	2.50	0.43
2:B:57:MET:CE	2:B:62:THR:HG22	2.48	0.43
2:D:233:LEU:HD23	2:D:233:LEU:HA	1.78	0.43
1:I:48:THR:HG22	1:I:51:ILE:CB	2.48	0.43
2:D:209:PRO:HD3	2:D:230:PHE:HE2	1.83	0.43
2:H:166:ILE:CG2	2:H:167:LYS:N	2.80	0.43
2:B:373:LEU:HD12	2:B:377:CYS:SG	2.58	0.43
2:L:116:SER:CB	2:L:142:THR:HG23	2.43	0.43
2:N:519:TRP:HA	2:N:568:ILE:O	2.18	0.43
1:K:137:PHE:HD1	2:L:17:VAL:CG2	2.31	0.43
1:A:27:GLN:HA	1:A:30:ALA:HB3	2.00	0.43
2:H:266:LEU:HD13	2:H:267:VAL:H	1.82	0.43
2:F:578:ARG:HG3	2:F:578:ARG:H	1.51	0.43
2:H:196:VAL:HG22	2:H:196:VAL:O	2.17	0.43
2:L:392:ASN:HD21	2:L:424:LEU:HA	1.84	0.43
2:H:318:PRO:C	2:H:320:LEU:H	2.22	0.43
1:M:26:SER:HB3	1:M:29:ILE:HG13	2.00	0.43
4:B:1100:OGK:C18	4:B:1100:OGK:N08	2.79	0.43
2:H:542:ILE:HD12	2:H:543:GLU:N	2.33	0.43
2:D:85:ARG:NH1	5:D:1101:PO4:O3	2.51	0.43
2:N:546:PRO:O	2:N:547:SER:CB	2.66	0.43
2:B:125:SER:O	2:B:129:LEU:HD22	2.19	0.43
2:P:398:ILE:C	2:P:400:THR:H	2.20	0.43
1:A:146:GLU:HA	2:D:539:TYR:OH	2.18	0.43
2:B:235:LEU:O	2:B:238:PHE:HB3	2.18	0.43
2:F:248:PHE:C	2:F:248:PHE:CD2	2.91	0.43
1:M:101:LEU:HD23	1:M:101:LEU:HA	1.80	0.43
2:P:289:PHE:N	2:P:290:PRO:CD	2.81	0.43
1:O:128:LYS:HB2	1:O:133:ILE:CD1	2.49	0.43
4:B:1100:OGK:N08	4:B:1100:OGK:H18A	2.24	0.43
2:B:519:TRP:C	2:B:519:TRP:CD1	2.90	0.43
2:N:289:PHE:N	2:N:290:PRO:CD	2.81	0.43
2:B:325:THR:O	2:B:349:ILE:HA	2.19	0.43
3:R:201:LEU:HD23	3:R:201:LEU:HA	1.85	0.43
1:C:113:LEU:O	1:C:117:THR:CG2	2.65	0.43
2:J:55:VAL:CG2	2:J:75:LEU:HD21	2.47	0.43
1:A:6:ILE:HG22	1:A:7:VAL:N	2.34	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:321:GLU:HA	2:F:344:LEU:HA	1.99	0.43
2:L:95:ASN:O	2:L:582:PRO:HG3	2.18	0.43
2:L:351:ARG:O	2:L:351:ARG:HG3	2.19	0.43
2:F:176:PHE:CZ	2:F:204:PHE:CZ	3.07	0.43
2:J:231:GLU:HG2	2:J:254:ASN:HD22	1.83	0.43
2:J:18:ASP:OD2	2:J:43:ARG:HD2	2.19	0.43
1:O:42:VAL:HG22	1:O:42:VAL:O	2.18	0.43
1:O:101:LEU:HD23	1:O:101:LEU:HA	1.88	0.43
2:P:343:GLN:NE2	2:P:343:GLN:HA	2.33	0.43
2:H:444:TYR:CD1	2:H:471:GLY:CA	3.01	0.43
2:J:308:GLU:O	2:J:312:THR:HG23	2.18	0.43
1:K:58:TYR:CE1	1:K:62:HIS:CE1	3.06	0.43
2:F:545:ILE:HG22	2:F:546:PRO:HD2	2.00	0.43
2:D:227:VAL:HG13	2:D:228:GLY:H	1.80	0.43
2:B:59:LEU:HD22	2:B:61:TYR:H	1.83	0.43
2:L:311:CYS:CB	2:L:336:VAL:HG21	2.48	0.43
2:L:398:ILE:C	2:L:400:THR:H	2.22	0.43
2:D:485:ARG:HE	2:D:485:ARG:HB3	1.69	0.43
2:L:152:THR:HG22	2:L:177:SER:HB2	2.01	0.43
2:N:304:LEU:HD12	2:N:304:LEU:HA	1.71	0.43
2:D:563:GLU:HG3	2:D:563:GLU:O	2.19	0.43
2:J:248:PHE:CD2	2:J:248:PHE:C	2.92	0.43
2:D:542:ILE:CD1	2:D:588:LEU:HD12	2.29	0.43
1:E:107:TYR:HE1	2:L:270:ARG:O	2.02	0.43
2:F:419:ILE:HD11	2:F:446:ARG:HH22	1.84	0.43
2:D:546:PRO:O	2:D:547:SER:CB	2.67	0.43
3:S:201:LEU:HD23	3:S:201:LEU:HA	1.79	0.43
2:B:157:LEU:HD12	2:B:157:LEU:O	2.19	0.43
2:H:367:GLN:HA	2:H:367:GLN:OE1	2.18	0.43
2:B:159:ILE:HG13	2:B:160:VAL:N	2.29	0.43
2:J:407:ASP:OD1	2:J:440:ARG:HD2	2.19	0.43
2:B:398:ILE:C	2:B:400:THR:H	2.22	0.43
2:H:301:LEU:HB3	2:H:302:TYR:CD2	2.53	0.43
2:H:207:ILE:H	2:H:207:ILE:HG12	1.59	0.43
1:C:8:LEU:HD23	1:C:42:VAL:CG1	2.49	0.43
2:J:266:LEU:HA	2:J:266:LEU:HD13	1.79	0.43
2:L:351:ARG:HD3	2:L:413:LEU:HD11	2.00	0.43
2:F:502:GLU:HG3	2:F:525:ALA:HA	2.00	0.43
1:M:42:VAL:HA	1:M:43:PRO:HD3	1.79	0.43
2:L:304:LEU:HD12	2:L:304:LEU:HA	1.75	0.43
2:B:116:SER:CB	2:B:142:THR:HG23	2.40	0.43
2:N:411:VAL:HG22	2:N:444:TYR:HB3	1.99	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:444:TYR:CE1	2:H:471:GLY:HA2	2.54	0.43
2:L:419:ILE:O	2:L:420:THR:C	2.56	0.43
2:D:519:TRP:HZ3	2:D:567:HIS:ND1	2.15	0.43
2:F:125:SER:O	2:F:128:ASP:OD2	2.37	0.43
2:D:59:LEU:O	2:D:62:THR:HB	2.18	0.43
2:N:490:LEU:HD23	2:N:490:LEU:HA	1.82	0.43
1:A:10:SER:OG	1:A:11:SER:N	2.52	0.43
2:J:59:LEU:HD22	2:J:61:TYR:H	1.84	0.43
2:H:129:LEU:HA	2:H:129:LEU:HD12	1.69	0.43
1:A:48:THR:HG23	1:A:51:ILE:H	1.84	0.43
1:M:148:GLU:HG2	1:M:152:ARG:HH12	1.83	0.43
2:H:480:LEU:HD22	2:H:500:PHE:CD1	2.53	0.43
2:P:46:LYS:HE3	2:P:46:LYS:HB2	1.54	0.43
2:N:412:LEU:HD12	2:N:412:LEU:O	2.18	0.43
2:J:542:ILE:CD1	2:J:588:LEU:HD12	2.29	0.43
2:B:411:VAL:HG22	2:B:444:TYR:HB3	2.01	0.43
2:H:325:THR:OG1	2:H:326:ARG:N	2.51	0.43
2:H:385:VAL:CG1	2:H:387:VAL:HG23	2.48	0.43
2:H:412:LEU:CD1	2:H:445:LEU:CD2	2.97	0.43
1:I:132:GLU:O	1:I:136:THR:HG23	2.19	0.43
1:A:108:LEU:HD12	1:A:110:ILE:CD1	2.42	0.43
2:N:472:TYR:CD2	2:N:497:GLY:O	2.71	0.43
2:F:136:ARG:HH21	2:L:164:ARG:HH21	1.65	0.43
2:H:298:LEU:HD22	2:H:300:LEU:CG	2.45	0.43
2:D:190:HIS:CD2	2:H:112:ARG:NH1	2.86	0.43
2:J:233:LEU:HA	2:J:233:LEU:HD23	1.81	0.43
1:M:134:ARG:HH11	2:N:40:VAL:HA	1.82	0.43
2:B:233:LEU:HD23	2:B:233:LEU:HA	1.80	0.43
3:U:201:LEU:HA	3:U:201:LEU:HD23	1.85	0.43
2:P:227:VAL:HG13	2:P:228:GLY:H	1.82	0.43
2:N:396:GLU:HG2	2:N:430:SER:OG	2.18	0.43
2:N:380:LEU:HA	2:N:380:LEU:HD23	1.77	0.43
2:F:65:PRO:HB3	2:F:103:TRP:CE3	2.54	0.43
1:M:153:ARG:CG	1:M:157:TRP:CZ3	3.02	0.43
2:B:201:MET:HB3	3:Q:211:ARG:HH12	1.82	0.43
5:B:1102:PO4:O3	3:Q:206:ARG:NH2	2.49	0.43
2:B:190:HIS:HE1	2:P:110:ASN:HA	1.84	0.43
1:E:91:MET:O	1:E:93:ILE:HG12	2.18	0.43
2:D:590:GLU:O	2:D:591:PRO:C	2.56	0.43
2:H:362:GLU:O	2:H:364:LEU:N	2.52	0.43
1:E:42:VAL:HA	1:E:43:PRO:HD3	1.79	0.43
1:E:128:LYS:HB2	1:E:133:ILE:CD1	2.49	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:565:PRO:HG2	3:V:201:LEU:HD21	1.99	0.43
2:F:311:CYS:CB	2:F:336:VAL:HG21	2.49	0.43
2:B:55:VAL:CG2	2:B:75:LEU:HD21	2.49	0.43
2:N:351:ARG:HD3	2:N:413:LEU:HD11	2.00	0.43
2:H:97:GLY:HA3	2:H:123:ILE:HD12	1.99	0.43
2:B:538:PRO:O	2:B:539:TYR:HB2	2.19	0.43
1:G:58:TYR:CD2	1:G:113:LEU:HD13	2.54	0.43
1:M:42:VAL:O	1:M:42:VAL:HG22	2.18	0.43
2:B:375:GLN:HG2	2:B:401:TYR:CE1	2.53	0.43
2:J:58:ALA:O	2:J:81:LYS:HB2	2.19	0.43
2:D:580:ASP:HA	2:N:206:LYS:HE3	2.01	0.43
2:H:316:LYS:C	2:H:318:PRO:HD3	2.40	0.42
2:H:450:LEU:HG	2:H:454:GLY:HA3	2.01	0.42
2:D:546:PRO:HB2	2:D:547:SER:H	1.72	0.42
2:F:519:TRP:CH2	2:F:567:HIS:CG	3.07	0.42
2:J:308:GLU:HG3	2:J:332:ARG:NH2	2.28	0.42
2:J:545:ILE:HG22	2:J:546:PRO:N	2.34	0.42
1:A:95:GLN:NE2	1:A:124:MET:CE	2.82	0.42
1:I:27:GLN:HA	1:I:30:ALA:HB3	2.01	0.42
2:P:578:ARG:HG3	2:P:578:ARG:H	1.44	0.42
2:D:263:TYR:O	2:D:266:LEU:HD23	2.19	0.42
2:L:231:GLU:HG2	2:L:254:ASN:HD22	1.83	0.42
1:I:42:VAL:HG22	1:I:42:VAL:O	2.18	0.42
2:H:437:LYS:HB3	2:H:437:LYS:HE2	1.80	0.42
2:D:343:GLN:NE2	2:D:343:GLN:HA	2.34	0.42
2:J:544:LEU:HD11	2:J:588:LEU:HD11	2.01	0.42
2:H:310:HIS:O	2:H:314:ILE:CG1	2.49	0.42
1:E:45:PRO:HD2	2:L:294:GLN:HB2	2.01	0.42
1:K:26:SER:C	1:K:28:THR:H	2.21	0.42
2:H:303:ALA:O	2:H:326:ARG:NH2	2.52	0.42
1:A:26:SER:HB3	1:A:29:ILE:HG13	2.01	0.42
2:B:523:TYR:CE2	2:B:568:ILE:HD11	2.54	0.42
1:E:26:SER:C	1:E:28:THR:H	2.22	0.42
2:H:304:LEU:HD22	2:H:304:LEU:HA	1.48	0.42
2:D:311:CYS:CB	2:D:336:VAL:HG21	2.50	0.42
2:B:344:LEU:HD23	2:B:380:LEU:HD21	2.01	0.42
2:P:338:ALA:O	2:P:376:GLY:HA3	2.19	0.42
2:B:198:ASN:OD1	2:B:198:ASN:C	2.57	0.42
2:B:46:LYS:HB2	2:B:46:LYS:HE3	1.51	0.42
2:J:289:PHE:N	2:J:290:PRO:CD	2.82	0.42
2:H:542:ILE:HD12	2:H:542:ILE:C	2.39	0.42
2:J:297:LYS:HE3	2:J:297:LYS:HB2	1.62	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:J:101:THR:N	2:J:102:PRO:CD	2.82	0.42
1:G:46:ASN:HB2	1:G:107:TYR:CZ	2.54	0.42
2:B:396:GLU:HG2	2:B:430:SER:OG	2.19	0.42
2:N:465:VAL:CG1	2:N:468:MET:HG3	2.48	0.42
1:C:27:GLN:HA	1:C:30:ALA:HB3	2.00	0.42
1:G:102:ILE:HG12	1:G:117:THR:OG1	2.18	0.42
2:H:20:VAL:O	2:H:24:VAL:HG23	2.20	0.42
2:D:380:LEU:HD23	2:D:380:LEU:HA	1.71	0.42
2:H:168:THR:CG2	2:H:196:VAL:HG13	2.50	0.42
2:L:299:ASP:OD2	2:L:301:LEU:HB2	2.18	0.42
2:N:412:LEU:HD12	2:N:412:LEU:C	2.40	0.42
2:D:78:LEU:HD12	2:D:79:LYS:H	1.85	0.42
2:D:231:GLU:HG2	2:D:254:ASN:HD22	1.84	0.42
2:F:198:ASN:OD1	2:F:198:ASN:C	2.56	0.42
2:N:544:LEU:HD11	2:N:588:LEU:CD1	2.50	0.42
1:E:46:ASN:HB2	1:E:107:TYR:CZ	2.54	0.42
2:J:367:GLN:CB	2:J:391:THR:HG22	2.37	0.42
2:H:350:GLU:CB	2:H:386:TYR:CD1	3.02	0.42
2:B:519:TRP:CH2	2:B:567:HIS:CE1	3.07	0.42
2:P:546:PRO:O	2:P:547:SER:CB	2.67	0.42
4:P:8100:OGK:N08	4:P:8100:OGK:H18A	2.25	0.42
4:D:2100:OGK:N08	4:D:2100:OGK:C18	2.81	0.42
2:H:208:SER:O	2:H:211:ASP:N	2.50	0.42
3:V:201:LEU:HA	3:V:202:PRO:HD3	1.82	0.42
2:N:280:MET:HG2	2:N:280:MET:O	2.19	0.42
2:B:398:ILE:HG23	2:B:402:LEU:HD11	2.02	0.42
2:N:477:ASP:N	2:N:477:ASP:OD1	2.52	0.42
1:M:87:ASP:HB3	1:M:116:LEU:CD2	2.49	0.42
1:A:148:GLU:HG2	1:A:152:ARG:HH12	1.84	0.42
2:N:232:ILE:HD12	2:N:252:SER:H	1.83	0.42
2:H:505:ILE:C	2:H:507:ALA:N	2.73	0.42
2:F:452:ASP:OD1	2:F:478:GLU:HB2	2.19	0.42
2:P:20:VAL:O	2:P:24:VAL:HG23	2.19	0.42
2:F:542:ILE:CG1	2:F:588:LEU:HB2	2.49	0.42
2:F:544:LEU:HD11	2:F:588:LEU:CD1	2.50	0.42
2:J:590:GLU:O	2:J:591:PRO:C	2.57	0.42
2:H:412:LEU:O	2:H:414:ASP:N	2.53	0.42
2:H:136:ARG:O	2:H:139:ASP:HB2	2.19	0.42
1:E:137:PHE:CD1	2:F:17:VAL:CG2	3.02	0.42
2:P:431:LEU:C	2:P:431:LEU:HD12	2.39	0.42
1:M:134:ARG:NE	1:M:141:ASN:HB2	2.34	0.42
1:K:48:THR:HG23	1:K:51:ILE:H	1.85	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:48:THR:HG23	1:I:51:ILE:H	1.85	0.42
2:L:221:SER:O	2:L:223:VAL:HG23	2.18	0.42
2:D:299:ASP:C	2:D:301:LEU:H	2.23	0.42
2:N:127:LEU:HD21	2:N:131:ARG:NH2	2.34	0.42
2:L:20:VAL:O	2:L:24:VAL:HG23	2.19	0.42
2:B:95:ASN:O	2:B:582:PRO:HG3	2.20	0.42
2:F:221:SER:O	2:F:223:VAL:HG23	2.20	0.42
2:J:563:GLU:O	2:J:563:GLU:HG3	2.19	0.42
2:H:362:GLU:HB3	2:H:363:GLY:H	1.52	0.42
2:L:519:TRP:C	2:L:519:TRP:CD1	2.90	0.42
2:H:518:LEU:HB3	2:H:570:ALA:HB3	2.02	0.42
2:H:194:LEU:HA	2:H:194:LEU:HD12	1.77	0.42
2:J:350:GLU:HB3	3:U:209:LEU:CD2	2.46	0.42
2:B:490:LEU:HD11	2:B:493:LEU:HD13	2.01	0.42
2:B:295:ILE:HG21	2:B:298:LEU:CD1	2.47	0.42
2:H:242:ALA:HB1	2:H:245:LEU:HB2	2.02	0.42
1:M:48:THR:HG23	1:M:51:ILE:H	1.84	0.42
2:F:227:VAL:HG13	2:F:228:GLY:H	1.84	0.42
2:D:125:SER:O	2:D:129:LEU:HD22	2.20	0.42
1:A:134:ARG:CZ	1:A:141:ASN:HD22	2.32	0.42
1:I:137:PHE:HD1	2:J:17:VAL:HG21	1.83	0.42
2:B:456:SER:CB	2:B:482:GLU:HB3	2.49	0.42
2:J:65:PRO:HA	2:J:103:TRP:CZ3	2.54	0.42
2:L:431:LEU:HD12	2:L:431:LEU:O	2.19	0.42
2:D:436:LYS:HE2	2:D:436:LYS:HB3	1.74	0.42
2:H:320:LEU:HD12	2:H:321:GLU:H	1.84	0.42
1:I:128:LYS:HB2	1:I:133:ILE:CD1	2.50	0.42
2:L:308:GLU:O	2:L:312:THR:HG23	2.20	0.42
1:C:132:GLU:O	1:C:136:THR:HG23	2.20	0.42
2:F:347:LEU:HD11	2:F:349:ILE:HD11	2.01	0.42
2:L:253:LEU:CD1	2:L:280:MET:HB2	2.46	0.42
2:J:101:THR:HG22	2:J:128:ASP:CB	2.48	0.42
2:P:59:LEU:O	2:P:62:THR:HB	2.19	0.42
1:I:101:LEU:HB3	1:I:117:THR:HG21	2.01	0.42
1:A:137:PHE:CD1	2:B:17:VAL:CG2	3.03	0.42
2:J:59:LEU:O	2:J:62:THR:HB	2.20	0.42
2:P:431:LEU:O	2:P:431:LEU:HD12	2.20	0.42
1:O:48:THR:HG22	1:O:51:ILE:HG12	2.01	0.42
1:A:98:LEU:CD2	1:A:120:THR:HG22	2.49	0.42
1:A:153:ARG:HG2	1:A:157:TRP:CZ3	2.54	0.42
2:H:115:LYS:O	2:H:141:GLU:N	2.51	0.42
2:F:339:GLN:HA	2:F:342:LYS:HE2	2.01	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:452:ASP:OD1	2:L:478:GLU:HB2	2.20	0.42
2:F:469:LEU:HA	2:F:494:GLU:O	2.19	0.42
2:P:231:GLU:HG2	2:P:254:ASN:HD22	1.84	0.42
2:N:299:ASP:OD2	2:N:301:LEU:HB2	2.19	0.42
2:B:299:ASP:OD2	2:B:301:LEU:HB2	2.19	0.42
2:F:563:GLU:HG3	2:F:563:GLU:O	2.20	0.42
1:G:91:MET:O	1:G:93:ILE:HG12	2.19	0.42
2:J:325:THR:OG1	2:J:326:ARG:N	2.53	0.42
2:H:535:MET:HE1	2:H:542:ILE:HG21	2.01	0.42
2:B:270:ARG:CB	1:O:107:TYR:HD1	2.31	0.42
1:M:10:SER:OG	1:M:11:SER:N	2.52	0.42
2:D:101:THR:HG22	2:D:128:ASP:CB	2.48	0.42
2:F:159:ILE:HD12	2:F:166:ILE:HD11	2.00	0.42
2:H:503:ARG:HE	2:H:503:ARG:H	1.66	0.42
2:F:78:LEU:HD12	2:F:79:LYS:H	1.84	0.42
2:F:211:ASP:O	2:F:215:ILE:HG13	2.19	0.42
1:G:101:LEU:HA	1:G:101:LEU:HD23	1.82	0.42
2:H:390:ILE:HG12	2:H:390:ILE:H	1.64	0.42
2:B:546:PRO:O	2:B:547:SER:CB	2.66	0.42
2:F:289:PHE:HD1	2:F:316:LYS:HD2	1.76	0.42
1:K:135:THR:HG22	1:K:136:THR:N	2.34	0.42
2:H:213:GLU:HA	2:H:238:PHE:HB2	2.01	0.42
2:J:477:ASP:N	2:J:477:ASP:OD1	2.53	0.42
2:N:429:ARG:HG2	2:N:433:ILE:HD12	2.02	0.42
2:L:412:LEU:HD12	2:L:412:LEU:O	2.19	0.42
1:O:59:CYS:O	1:O:63:VAL:HG23	2.19	0.42
1:A:8:LEU:HD23	1:A:42:VAL:CG1	2.50	0.42
2:P:563:GLU:HG3	2:P:563:GLU:O	2.20	0.42
2:H:285:MET:O	2:H:287:ILE:N	2.53	0.42
2:N:419:ILE:O	2:N:420:THR:O	2.37	0.42
1:M:99:PHE:CE2	2:N:17:VAL:N	2.88	0.42
2:H:462:SER:HB2	2:H:465:VAL:HB	2.02	0.42
2:F:380:LEU:HD23	2:F:380:LEU:HA	1.74	0.42
2:P:392:ASN:ND2	2:P:421:ASP:OD1	2.48	0.42
2:D:431:LEU:HD12	2:D:431:LEU:C	2.39	0.42
2:D:419:ILE:O	2:D:420:THR:O	2.38	0.41
2:L:546:PRO:O	2:L:547:SER:CB	2.67	0.41
2:B:300:LEU:HD12	2:B:323:LEU:HD11	2.02	0.41
2:F:314:ILE:HD11	2:F:329:ILE:HD11	2.01	0.41
1:M:134:ARG:NH1	2:N:40:VAL:HA	2.35	0.41
2:D:396:GLU:HG2	2:D:430:SER:OG	2.20	0.41
2:L:396:GLU:HG2	2:L:430:SER:OG	2.19	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:188:ALA:HB1	2:H:214:THR:OG1	2.20	0.41
2:B:272:LEU:C	2:B:272:LEU:CD2	2.88	0.41
2:D:296:ARG:HH12	1:G:32:MET:HG3	1.85	0.41
2:D:543:GLU:OE2	2:D:578:ARG:HD3	2.20	0.41
2:N:326:ARG:HA	2:N:350:GLU:O	2.20	0.41
1:A:102:ILE:HG12	1:A:117:THR:HB	2.02	0.41
2:H:100:VAL:O	2:H:100:VAL:HG13	2.19	0.41
2:L:285:MET:N	2:L:286:PRO:CD	2.83	0.41
2:F:282:PRO:HA	2:F:285:MET:CE	2.48	0.41
2:F:285:MET:N	2:F:286:PRO:CD	2.81	0.41
2:L:308:GLU:HG3	2:L:332:ARG:NH2	2.26	0.41
2:P:253:LEU:CD1	2:P:280:MET:HB2	2.49	0.41
1:I:101:LEU:HD23	1:I:101:LEU:HA	1.83	0.41
2:H:428:VAL:CG2	2:H:429:ARG:N	2.83	0.41
1:A:48:THR:HG22	1:A:51:ILE:HG12	2.03	0.41
2:J:143:LEU:CD2	2:J:159:ILE:HD13	2.50	0.41
3:Q:201:LEU:HA	3:Q:201:LEU:HD23	1.85	0.41
2:H:184:LEU:HD23	2:H:184:LEU:HA	1.69	0.41
2:J:314:ILE:HD11	2:J:329:ILE:HD11	2.01	0.41
2:B:223:VAL:O	2:B:245:LEU:HD12	2.20	0.41
2:D:337:LEU:HD12	2:D:341:CYS:SG	2.60	0.41
2:N:544:LEU:HD11	2:N:588:LEU:HD11	2.01	0.41
2:H:87:ALA:HB2	2:H:92:ILE:CB	2.50	0.41
1:I:129:THR:HB	1:I:130:PRO:HD2	2.03	0.41
1:O:125:ILE:HG23	1:O:133:ILE:CD1	2.44	0.41
1:I:106:ASN:HB2	1:I:114:LEU:HD11	2.02	0.41
1:E:134:ARG:NE	1:E:141:ASN:HB2	2.35	0.41
2:J:322:VAL:HG13	2:J:346:ARG:HB2	2.01	0.41
1:O:105:ALA:HB2	1:O:113:LEU:HD23	2.02	0.41
2:J:75:LEU:HD23	2:J:75:LEU:HA	1.83	0.41
1:M:27:GLN:HA	1:M:30:ALA:HB3	2.02	0.41
1:K:46:ASN:HB2	1:K:107:TYR:CZ	2.55	0.41
2:N:502:GLU:HG3	2:N:525:ALA:HA	2.03	0.41
2:N:358:MET:HG3	2:N:364:LEU:HG	2.01	0.41
2:L:209:PRO:HD3	2:L:230:PHE:HE2	1.84	0.41
2:B:348:ARG:HG3	2:B:384:ALA:HB3	2.02	0.41
2:N:107:ILE:HA	2:N:111:LEU:HB2	2.02	0.41
1:I:19:GLU:OE1	1:I:19:GLU:N	2.51	0.41
2:B:563:GLU:HG3	2:B:563:GLU:O	2.21	0.41
2:N:590:GLU:O	2:N:591:PRO:C	2.57	0.41
2:L:289:PHE:N	2:L:290:PRO:CD	2.82	0.41
2:D:325:THR:O	2:D:349:ILE:HA	2.20	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:347:LEU:CD2	2:H:373:LEU:HD21	2.50	0.41
2:P:83:LYS:HA	2:P:84:PRO:HD3	1.86	0.41
1:C:105:ALA:HB2	1:C:113:LEU:HD23	2.02	0.41
2:L:125:SER:O	2:L:129:LEU:HD22	2.20	0.41
2:H:395:LEU:HB3	2:H:427:GLY:C	2.40	0.41
2:F:343:GLN:NE2	2:F:343:GLN:HA	2.35	0.41
2:H:411:VAL:HG23	4:H:4100:OGK:H16A	2.02	0.41
1:M:125:ILE:HG23	1:M:133:ILE:CD1	2.39	0.41
2:N:519:TRP:HZ3	2:N:567:HIS:ND1	2.18	0.41
2:P:328:VAL:CG2	2:P:359:GLU:HB2	2.43	0.41
2:L:289:PHE:HB2	2:L:290:PRO:HD3	2.01	0.41
2:P:308:GLU:O	2:P:312:THR:HG23	2.21	0.41
1:A:132:GLU:O	1:A:136:THR:HG23	2.20	0.41
2:P:490:LEU:HD11	2:P:493:LEU:HD13	2.03	0.41
2:H:124:VAL:CG1	2:H:129:LEU:HD13	2.51	0.41
1:K:137:PHE:CD1	2:L:17:VAL:CG2	3.03	0.41
1:O:102:ILE:HD12	2:P:20:VAL:HG21	2.02	0.41
2:P:89:PHE:CD1	3:X:206:ARG:HA	2.56	0.41
2:F:54:HIS:HE1	2:F:56:THR:OG1	2.03	0.41
2:B:266:LEU:HD13	2:B:266:LEU:HA	1.83	0.41
2:D:187:LEU:HD23	2:D:187:LEU:HA	1.88	0.41
2:P:436:LYS:HB3	2:P:436:LYS:HE2	1.74	0.41
2:L:590:GLU:O	2:L:591:PRO:C	2.57	0.41
2:H:365:VAL:HG21	2:H:387:VAL:CB	2.50	0.41
2:D:285:MET:N	2:D:286:PRO:HD2	2.36	0.41
2:L:519:TRP:CH2	2:L:567:HIS:CE1	3.08	0.41
1:A:114:LEU:HD12	1:A:114:LEU:HA	1.97	0.41
2:F:311:CYS:SG	2:F:333:GLY:HA2	2.60	0.41
1:O:153:ARG:NH2	2:P:539:TYR:HE1	2.19	0.41
2:H:54:HIS:O	2:H:55:VAL:HG23	2.20	0.41
2:H:111:LEU:C	2:H:113:GLN:N	2.74	0.41
1:A:101:LEU:HB3	1:A:117:THR:HG21	2.01	0.41
2:P:545:ILE:HB	2:P:567:HIS:HB2	2.01	0.41
2:J:409:ARG:HD3	4:J:5100:OGK:H13A	2.01	0.41
2:N:308:GLU:HG3	2:N:332:ARG:NH2	2.30	0.41
1:I:134:ARG:NE	1:I:141:ASN:HB2	2.36	0.41
2:H:315:GLN:HG3	2:H:340:TYR:CZ	2.55	0.41
1:K:8:LEU:HD23	1:K:42:VAL:CG1	2.50	0.41
2:L:431:LEU:HD12	2:L:431:LEU:C	2.41	0.41
2:B:231:GLU:HG2	2:B:254:ASN:HD22	1.86	0.41
2:J:209:PRO:HD3	2:J:230:PHE:HE2	1.85	0.41
2:N:586:ARG:NH2	2:N:588:LEU:HD21	2.35	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:52:ARG:HH11	2:H:72:PHE:HE2	1.69	0.41
2:P:590:GLU:O	2:P:591:PRO:C	2.58	0.41
2:H:408:PHE:C	2:H:409:ARG:HG2	2.41	0.41
2:N:496:ARG:HA	2:N:521:GLN:O	2.19	0.41
2:P:490:LEU:HA	2:P:490:LEU:HD23	1.72	0.41
2:J:124:VAL:CG1	2:J:129:LEU:HD13	2.50	0.41
2:L:101:THR:N	2:L:102:PRO:CD	2.84	0.41
1:C:137:PHE:HD1	2:D:17:VAL:CG2	2.34	0.41
2:H:168:THR:HG22	2:H:196:VAL:HG13	2.01	0.41
2:N:121:ARG:NH2	5:N:1103:PO4:O4	2.54	0.41
2:D:429:ARG:HG2	2:D:433:ILE:HD12	2.02	0.41
2:P:198:ASN:OD1	2:P:198:ASN:C	2.59	0.41
2:D:405:LEU:HA	2:D:405:LEU:HD23	1.88	0.41
2:L:533:MET:CE	2:L:588:LEU:HD13	2.51	0.41
2:H:455:LEU:HD21	2:H:473:VAL:HG21	2.03	0.41
2:H:92:ILE:HD11	2:H:519:TRP:CZ3	2.56	0.41
2:B:328:VAL:CG2	2:B:359:GLU:HB2	2.45	0.41
2:H:468:MET:HE2	2:H:470:LEU:HD21	2.02	0.41
2:B:519:TRP:HA	2:B:568:ILE:O	2.21	0.41
2:D:545:ILE:HG22	2:D:546:PRO:CD	2.51	0.41
2:H:170:LEU:HA	2:H:198:ASN:O	2.20	0.41
1:O:46:ASN:HB2	1:O:107:TYR:CZ	2.56	0.41
2:J:545:ILE:HB	2:J:567:HIS:HB2	2.03	0.41
2:N:308:GLU:O	2:N:312:THR:HG23	2.21	0.41
2:H:391:THR:O	2:H:393:GLU:N	2.54	0.41
2:H:59:LEU:O	2:H:62:THR:HB	2.21	0.41
2:L:490:LEU:HD23	2:L:490:LEU:HA	1.77	0.41
2:H:584:THR:HG22	2:H:584:THR:O	2.21	0.41
2:J:191:ASN:ND2	2:J:192:THR:N	2.68	0.41
2:N:398:ILE:HG22	2:N:431:LEU:HD13	2.02	0.41
2:B:314:ILE:HD11	2:B:329:ILE:HD11	2.03	0.41
1:G:129:THR:OG1	1:G:132:GLU:HG3	2.21	0.41
1:C:99:PHE:CE2	2:D:17:VAL:HG22	2.56	0.41
1:G:148:GLU:HG3	1:G:152:ARG:HH11	1.85	0.41
2:P:392:ASN:HD21	2:P:424:LEU:HA	1.85	0.41
2:B:78:LEU:HD12	2:B:79:LYS:H	1.86	0.41
2:J:405:LEU:HA	2:J:405:LEU:HD23	1.85	0.41
2:F:338:ALA:O	2:F:376:GLY:HA3	2.21	0.41
2:J:412:LEU:C	2:J:412:LEU:HD12	2.41	0.41
2:H:70:ARG:O	2:H:72:PHE:N	2.54	0.41
2:J:367:GLN:NE2	2:J:389:ASP:OD1	2.54	0.41
1:A:26:SER:C	1:A:28:THR:H	2.24	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:J:328:VAL:CG2	2:J:359:GLU:HB2	2.45	0.41
1:C:125:ILE:HG23	1:C:133:ILE:CD1	2.44	0.41
2:H:98:GLY:C	2:H:122:MET:HE3	2.41	0.41
2:J:519:TRP:C	2:J:519:TRP:CD1	2.92	0.41
1:O:114:LEU:HD12	1:O:114:LEU:HA	1.91	0.41
1:A:141:ASN:C	1:A:141:ASN:OD1	2.58	0.41
2:H:352:GLY:O	2:H:353:ALA:CB	2.69	0.41
2:F:480:LEU:HB2	2:F:500:PHE:CE2	2.56	0.41
2:L:436:LYS:HE2	2:L:436:LYS:HB3	1.70	0.41
1:G:147:GLU:HG3	1:G:147:GLU:H	1.60	0.41
2:D:489:ASN:HA	2:D:489:ASN:HD22	1.65	0.41
2:J:533:MET:C	2:J:535:MET:H	2.24	0.40
2:D:533:MET:C	2:D:535:MET:H	2.25	0.40
2:B:533:MET:CE	2:B:588:LEU:HD13	2.47	0.40
1:G:130:PRO:HD3	2:H:36:SER:CB	2.50	0.40
1:I:26:SER:C	1:I:28:THR:H	2.25	0.40
2:L:409:ARG:HD3	4:L:6100:OGK:H13A	2.03	0.40
2:H:120:ARG:O	2:H:122:MET:HG3	2.20	0.40
2:P:101:THR:N	2:P:102:PRO:CD	2.84	0.40
1:C:158:ALA:HA	2:D:62:THR:HG23	2.03	0.40
1:C:10:SER:OG	1:C:11:SER:N	2.54	0.40
2:P:399:GLY:HA3	2:P:431:LEU:HA	2.04	0.40
2:D:404:ASN:HB3	2:D:437:LYS:HD2	2.03	0.40
2:F:263:TYR:O	2:F:266:LEU:HD23	2.20	0.40
2:B:152:THR:HG22	2:B:177:SER:HB2	2.03	0.40
1:K:114:LEU:C	1:K:116:LEU:N	2.73	0.40
2:P:276:GLY:HA3	2:P:299:ASP:HB3	2.02	0.40
2:P:266:LEU:HA	2:P:266:LEU:HD13	1.86	0.40
2:L:327:ASN:HD22	2:L:327:ASN:N	2.18	0.40
2:L:266:LEU:HA	2:L:266:LEU:HD13	1.83	0.40
2:P:305:LEU:HD23	2:P:305:LEU:O	2.19	0.40
2:N:563:GLU:HG3	2:N:563:GLU:O	2.21	0.40
3:W:201:LEU:HD23	3:W:201:LEU:HA	1.81	0.40
2:N:285:MET:N	2:N:286:PRO:CD	2.85	0.40
1:E:130:PRO:O	1:E:134:ARG:HG2	2.20	0.40
2:N:55:VAL:HG21	2:N:72:PHE:CD1	2.56	0.40
2:N:545:ILE:HG22	2:N:546:PRO:N	2.36	0.40
2:D:490:LEU:HD23	2:D:490:LEU:HA	1.74	0.40
2:F:55:VAL:HG21	2:F:72:PHE:CD1	2.56	0.40
2:P:232:ILE:O	2:P:235:LEU:HB2	2.21	0.40
2:H:274:ARG:HA	2:H:297:LYS:HB3	2.03	0.40
2:F:392:ASN:ND2	2:F:421:ASP:OD1	2.54	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:429:ARG:HG2	2:F:433:ILE:HD12	2.04	0.40
2:B:127:LEU:HB2	2:F:126:ASP:HB3	2.03	0.40
2:P:95:ASN:O	2:P:582:PRO:HG3	2.21	0.40
2:B:489:ASN:HA	2:B:489:ASN:HD22	1.66	0.40
2:B:304:LEU:HD12	2:B:304:LEU:HA	1.72	0.40
2:H:32:LYS:HD3	2:H:32:LYS:HA	1.81	0.40
2:F:118:HIS:CD2	2:F:118:HIS:C	2.94	0.40
2:J:436:LYS:HE2	2:J:436:LYS:HB3	1.77	0.40
2:N:328:VAL:CG2	2:N:359:GLU:HB2	2.41	0.40
2:B:545:ILE:HG22	2:B:546:PRO:N	2.36	0.40
2:F:289:PHE:N	2:F:290:PRO:CD	2.84	0.40
2:P:519:TRP:HZ3	2:P:567:HIS:ND1	2.20	0.40
2:L:519:TRP:HZ3	2:L:567:HIS:ND1	2.20	0.40
2:B:83:LYS:HA	2:B:84:PRO:HD3	1.86	0.40
2:P:184:LEU:HA	2:P:184:LEU:HD23	1.85	0.40
2:F:398:ILE:C	2:F:400:THR:H	2.25	0.40
2:P:351:ARG:HD3	2:P:413:LEU:HD11	2.03	0.40
1:I:153:ARG:CG	1:I:157:TRP:CZ3	3.04	0.40
1:I:146:GLU:HA	2:L:539:TYR:OH	2.22	0.40
2:N:501:SER:O	2:N:505:ILE:HG12	2.21	0.40
2:B:263:TYR:O	2:B:266:LEU:HD23	2.21	0.40
2:P:107:ILE:HA	2:P:111:LEU:HB2	2.03	0.40
2:P:100:VAL:HG11	2:P:124:VAL:HG22	2.03	0.40
2:P:187:LEU:HA	2:P:187:LEU:HD23	1.85	0.40
2:P:327:ASN:N	2:P:327:ASN:HD22	2.20	0.40
1:K:26:SER:CB	1:K:29:ILE:HG13	2.50	0.40
4:H:4100:OGK:N08	4:H:4100:OGK:C18	2.84	0.40
1:O:132:GLU:O	1:O:136:THR:HG23	2.22	0.40
1:E:26:SER:CB	1:E:29:ILE:HG13	2.51	0.40
2:F:308:GLU:HG3	2:F:332:ARG:NH2	2.28	0.40
2:H:191:ASN:ND2	2:H:193:SER:N	2.70	0.40
1:K:114:LEU:C	1:K:116:LEU:H	2.25	0.40
2:P:263:TYR:O	2:P:266:LEU:HD23	2.21	0.40
1:G:135:THR:HG22	1:G:136:THR:HG23	2.04	0.40
2:N:124:VAL:O	2:N:151:PHE:HB3	2.20	0.40
2:B:319:ASN:ND2	1:O:43:PRO:HG2	2.36	0.40
1:O:148:GLU:HG2	1:O:152:ARG:HH12	1.85	0.40
2:N:538:PRO:O	2:N:539:TYR:HB2	2.21	0.40
2:L:343:GLN:NE2	2:L:343:GLN:HA	2.36	0.40
2:P:289:PHE:N	2:P:290:PRO:HD2	2.37	0.40
2:B:285:MET:N	2:B:286:PRO:CD	2.85	0.40
2:N:492:LYS:HG3	2:N:517:TYR:CD2	2.57	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:N:289:PHE:N	2:N:290:PRO:HD2	2.37	0.40
2:F:519:TRP:HZ3	2:F:567:HIS:ND1	2.17	0.40
2:H:80:LEU:CD1	2:H:103:TRP:CG	3.05	0.40
2:H:191:ASN:C	2:H:191:ASN:HD22	2.24	0.40
2:H:374:ALA:CB	2:H:398:ILE:HD12	2.49	0.40
2:J:253:LEU:CD1	2:J:280:MET:HB2	2.49	0.40
2:F:398:ILE:HG22	2:F:431:LEU:HD13	2.03	0.40
2:P:400:THR:O	2:P:403:LYS:HE2	2.22	0.40
2:N:398:ILE:CG2	2:N:431:LEU:HD13	2.52	0.40
2:J:375:GLN:HG2	2:J:401:TYR:CE1	2.56	0.40
2:H:453:LEU:O	2:H:456:SER:HB3	2.21	0.40
1:E:148:GLU:HG2	1:E:152:ARG:HH12	1.86	0.40
2:P:358:MET:HG3	2:P:364:LEU:HG	2.03	0.40
1:O:61:ARG:HA	1:O:61:ARG:HD2	1.97	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	140/160 (88%)	115 (82%)	20 (14%)	5 (4%)	5	42
1	C	140/160 (88%)	116 (83%)	19 (14%)	5 (4%)	5	42
1	E	140/160 (88%)	118 (84%)	17 (12%)	5 (4%)	5	42
1	G	140/160 (88%)	113 (81%)	21 (15%)	6 (4%)	4	35
1	I	140/160 (88%)	117 (84%)	17 (12%)	6 (4%)	4	35
1	K	140/160 (88%)	116 (83%)	19 (14%)	5 (4%)	5	42
1	M	140/160 (88%)	115 (82%)	20 (14%)	5 (4%)	5	42
1	O	140/160 (88%)	115 (82%)	19 (14%)	6 (4%)	4	35
2	B	564/592 (95%)	501 (89%)	57 (10%)	6 (1%)	21	72
2	D	564/592 (95%)	503 (89%)	53 (9%)	8 (1%)	16	66
2	F	564/592 (95%)	505 (90%)	52 (9%)	7 (1%)	19	70

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	558/592 (94%)	426 (76%)	95 (17%)	37 (7%)	2	21
2	J	564/592 (95%)	505 (90%)	52 (9%)	7 (1%)	19	70
2	L	564/592 (95%)	501 (89%)	55 (10%)	8 (1%)	16	66
2	N	564/592 (95%)	502 (89%)	55 (10%)	7 (1%)	19	70
2	P	564/592 (95%)	500 (89%)	58 (10%)	6 (1%)	21	72
3	Q	16/21 (76%)	15 (94%)	1 (6%)	0	100	100
3	R	16/21 (76%)	14 (88%)	2 (12%)	0	100	100
3	S	16/21 (76%)	14 (88%)	2 (12%)	0	100	100
3	U	16/21 (76%)	14 (88%)	2 (12%)	0	100	100
3	V	16/21 (76%)	14 (88%)	2 (12%)	0	100	100
3	W	16/21 (76%)	15 (94%)	1 (6%)	0	100	100
3	X	16/21 (76%)	14 (88%)	2 (12%)	0	100	100
All	All	5738/6163 (93%)	4968 (87%)	641 (11%)	129 (2%)	10	56

All (129) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	92	LYS
2	B	420	THR
2	B	546	PRO
2	D	420	THR
2	D	546	PRO
2	F	420	THR
2	F	546	PRO
2	H	41	CYS
2	H	71	ARG
2	H	172	GLU
2	H	200	TYR
2	H	270	ARG
2	H	278	SER
2	H	357	GLY
2	H	364	LEU
2	H	425	ASP
2	H	525	ALA
2	H	526	SER
2	J	420	THR
2	J	546	PRO
1	K	92	LYS

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Mol	Chain	Res	Type
2	L	420	THR
2	L	546	PRO
2	N	420	THR
2	N	546	PRO
2	P	420	THR
2	P	546	PRO
1	A	10	SER
1	A	13	GLY
1	A	36	ASP
1	A	126	LYS
2	B	547	SER
1	C	10	SER
1	C	13	GLY
1	C	36	ASP
1	C	92	LYS
1	C	126	LYS
2	D	547	SER
1	E	10	SER
1	E	13	GLY
1	E	36	ASP
1	E	92	LYS
1	E	126	LYS
1	G	10	SER
1	G	13	GLY
1	G	36	ASP
1	G	92	LYS
1	G	128	LYS
2	H	97	GLY
2	H	228	GLY
2	H	271	LYS
2	H	279	TYR
2	H	353	ALA
2	H	365	VAL
2	H	392	ASN
2	H	449	GLY
2	H	528	THR
1	I	10	SER
1	I	13	GLY
1	I	36	ASP
1	I	92	LYS
1	I	126	LYS
2	J	547	SER

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Mol	Chain	Res	Type
1	K	10	SER
1	K	36	ASP
1	K	126	LYS
2	L	547	SER
1	M	10	SER
1	M	13	GLY
1	M	36	ASP
1	M	92	LYS
1	M	126	LYS
1	O	10	SER
1	O	36	ASP
1	O	92	LYS
1	O	126	LYS
2	P	547	SER
2	B	590	GLU
2	D	590	GLU
2	F	20	VAL
2	F	547	SER
2	F	590	GLU
2	H	70	ARG
2	H	319	ASN
2	H	499	CYS
2	H	533	MET
2	H	577	GLN
2	H	582	PRO
2	J	590	GLU
1	K	13	GLY
2	L	590	GLU
2	N	547	SER
2	N	590	GLU
1	O	13	GLY
2	P	590	GLU
2	B	20	VAL
2	D	20	VAL
2	D	112	ARG
2	H	112	ARG
2	H	137	ALA
2	H	269	PRO
2	H	305	LEU
2	H	358	MET
2	H	538	PRO
2	L	112	ARG

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Mol	Chain	Res	Type
2	N	290	PRO
2	P	20	VAL
2	D	290	PRO
2	H	206	LYS
1	I	137	PHE
2	J	20	VAL
2	J	290	PRO
2	L	20	VAL
2	L	290	PRO
2	N	20	VAL
1	O	27	GLN
2	P	290	PRO
2	F	290	PRO
1	G	153	ARG
2	H	209	PRO
2	H	276	GLY
2	H	20	VAL
2	H	286	PRO
2	B	591	PRO
2	D	591	PRO
2	N	591	PRO
2	F	591	PRO
2	J	591	PRO
2	L	591	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	127/137 (93%)	114 (90%)	13 (10%)	11	42
1	C	127/137 (93%)	115 (91%)	12 (9%)	13	47
1	E	127/137 (93%)	113 (89%)	14 (11%)	9	38
1	G	127/137 (93%)	113 (89%)	14 (11%)	9	38
1	I	127/137 (93%)	114 (90%)	13 (10%)	11	42
1	K	127/137 (93%)	114 (90%)	13 (10%)	11	42
1	M	127/137 (93%)	114 (90%)	13 (10%)	11	42

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	O	127/137 (93%)	113 (89%)	14 (11%)	9	38
2	B	500/523 (96%)	433 (87%)	67 (13%)	6	27
2	D	500/523 (96%)	432 (86%)	68 (14%)	5	26
2	F	500/523 (96%)	433 (87%)	67 (13%)	6	27
2	H	494/523 (94%)	397 (80%)	97 (20%)	2	9
2	J	500/523 (96%)	433 (87%)	67 (13%)	6	27
2	L	500/523 (96%)	431 (86%)	69 (14%)	5	25
2	N	500/523 (96%)	432 (86%)	68 (14%)	5	26
2	P	500/523 (96%)	434 (87%)	66 (13%)	6	28
3	Q	16/19 (84%)	16 (100%)	0	100	100
3	R	16/19 (84%)	16 (100%)	0	100	100
3	S	16/19 (84%)	16 (100%)	0	100	100
3	U	16/19 (84%)	16 (100%)	0	100	100
3	V	16/19 (84%)	16 (100%)	0	100	100
3	W	16/19 (84%)	16 (100%)	0	100	100
3	X	16/19 (84%)	16 (100%)	0	100	100
All	All	5122/5413 (95%)	4447 (87%)	675 (13%)	6	28

All (675) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LEU
1	A	14	GLU
1	A	22	VAL
1	A	29	ILE
1	A	35	ASP
1	A	36	ASP
1	A	42	VAL
1	A	44	LEU
1	A	48	THR
1	A	81	ASP
1	A	125	ILE
1	A	153	ARG
1	A	160	GLU
2	B	16	THR
2	B	20	VAL

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Mol	Chain	Res	Type
2	B	35	ASP
2	B	40	VAL
2	B	43	ARG
2	B	48	ASP
2	B	52	ARG
2	B	59	LEU
2	B	62	THR
2	B	64	THR
2	B	104	VAL
2	B	123	ILE
2	B	129	LEU
2	B	132	LEU
2	B	136	ARG
2	B	142	THR
2	B	159	ILE
2	B	166	ILE
2	B	168	THR
2	B	182	LYS
2	B	184	LEU
2	B	191	ASN
2	B	193	SER
2	B	196	VAL
2	B	214	THR
2	B	220	ARG
2	B	225	VAL
2	B	230	PHE
2	B	231	GLU
2	B	232	ILE
2	B	270	ARG
2	B	272	LEU
2	B	275	LEU
2	B	284	GLU
2	B	301	LEU
2	B	304	LEU
2	B	305	LEU
2	B	312	THR
2	B	314	ILE
2	B	327	ASN
2	B	349	ILE
2	B	360	ASP
2	B	367	GLN
2	B	377	CYS

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Mol	Chain	Res	Type
2	B	390	ILE
2	B	400	THR
2	B	402	LEU
2	B	404	ASN
2	B	405	LEU
2	B	412	LEU
2	B	418	ARG
2	B	422	LEU
2	B	447	GLN
2	B	453	LEU
2	B	455	LEU
2	B	477	ASP
2	B	481	MET
2	B	490	LEU
2	B	495	MET
2	B	503	ARG
2	B	510	THR
2	B	516	ARG
2	B	519	TRP
2	B	528	THR
2	B	537	ARG
2	B	542	ILE
2	B	578	ARG
1	C	8	LEU
1	C	14	GLU
1	C	29	ILE
1	C	35	ASP
1	C	36	ASP
1	C	42	VAL
1	C	44	LEU
1	C	48	THR
1	C	125	ILE
1	C	150	GLU
1	C	153	ARG
1	C	160	GLU
2	D	16	THR
2	D	20	VAL
2	D	35	ASP
2	D	40	VAL
2	D	43	ARG
2	D	48	ASP
2	D	52	ARG

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Mol	Chain	Res	Type
2	D	59	LEU
2	D	62	THR
2	D	64	THR
2	D	104	VAL
2	D	123	ILE
2	D	129	LEU
2	D	132	LEU
2	D	136	ARG
2	D	142	THR
2	D	159	ILE
2	D	166	ILE
2	D	168	THR
2	D	182	LYS
2	D	184	LEU
2	D	191	ASN
2	D	196	VAL
2	D	214	THR
2	D	220	ARG
2	D	225	VAL
2	D	230	PHE
2	D	231	GLU
2	D	232	ILE
2	D	270	ARG
2	D	272	LEU
2	D	275	LEU
2	D	284	GLU
2	D	301	LEU
2	D	304	LEU
2	D	305	LEU
2	D	312	THR
2	D	314	ILE
2	D	327	ASN
2	D	349	ILE
2	D	360	ASP
2	D	367	GLN
2	D	377	CYS
2	D	390	ILE
2	D	400	THR
2	D	402	LEU
2	D	404	ASN
2	D	405	LEU
2	D	412	LEU

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Mol	Chain	Res	Type
2	D	413	LEU
2	D	418	ARG
2	D	422	LEU
2	D	447	GLN
2	D	453	LEU
2	D	455	LEU
2	D	477	ASP
2	D	481	MET
2	D	490	LEU
2	D	495	MET
2	D	496	ARG
2	D	503	ARG
2	D	510	THR
2	D	516	ARG
2	D	519	TRP
2	D	528	THR
2	D	537	ARG
2	D	542	ILE
2	D	578	ARG
1	E	8	LEU
1	E	14	GLU
1	E	22	VAL
1	E	29	ILE
1	E	35	ASP
1	E	36	ASP
1	E	42	VAL
1	E	44	LEU
1	E	48	THR
1	E	81	ASP
1	E	125	ILE
1	E	150	GLU
1	E	153	ARG
1	E	160	GLU
2	F	16	THR
2	F	20	VAL
2	F	35	ASP
2	F	40	VAL
2	F	43	ARG
2	F	48	ASP
2	F	52	ARG
2	F	59	LEU
2	F	62	THR

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Mol	Chain	Res	Type
2	F	64	THR
2	F	104	VAL
2	F	123	ILE
2	F	129	LEU
2	F	132	LEU
2	F	136	ARG
2	F	142	THR
2	F	159	ILE
2	F	166	ILE
2	F	168	THR
2	F	182	LYS
2	F	184	LEU
2	F	191	ASN
2	F	196	VAL
2	F	201	MET
2	F	214	THR
2	F	220	ARG
2	F	225	VAL
2	F	231	GLU
2	F	232	ILE
2	F	270	ARG
2	F	272	LEU
2	F	275	LEU
2	F	284	GLU
2	F	301	LEU
2	F	304	LEU
2	F	305	LEU
2	F	312	THR
2	F	314	ILE
2	F	327	ASN
2	F	349	ILE
2	F	360	ASP
2	F	367	GLN
2	F	377	CYS
2	F	390	ILE
2	F	400	THR
2	F	402	LEU
2	F	404	ASN
2	F	405	LEU
2	F	412	LEU
2	F	413	LEU
2	F	418	ARG

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Mol	Chain	Res	Type
2	F	422	LEU
2	F	447	GLN
2	F	453	LEU
2	F	455	LEU
2	F	477	ASP
2	F	481	MET
2	F	490	LEU
2	F	495	MET
2	F	503	ARG
2	F	510	THR
2	F	516	ARG
2	F	519	TRP
2	F	528	THR
2	F	537	ARG
2	F	542	ILE
2	F	578	ARG
1	G	8	LEU
1	G	14	GLU
1	G	22	VAL
1	G	35	ASP
1	G	36	ASP
1	G	42	VAL
1	G	44	LEU
1	G	48	THR
1	G	81	ASP
1	G	129	THR
1	G	133	ILE
1	G	144	THR
1	G	150	GLU
1	G	153	ARG
2	H	14	VAL
2	H	17	VAL
2	H	20	VAL
2	H	21	ILE
2	H	35	ASP
2	H	40	VAL
2	H	46	LYS
2	H	52	ARG
2	H	59	LEU
2	H	62	THR
2	H	64	THR
2	H	71	ARG

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Mol	Chain	Res	Type
2	H	75	LEU
2	H	76	ARG
2	H	77	SER
2	H	85	ARG
2	H	95	ASN
2	H	100	VAL
2	H	111	LEU
2	H	128	ASP
2	H	132	LEU
2	H	136	ARG
2	H	140	LEU
2	H	142	THR
2	H	159	ILE
2	H	166	ILE
2	H	168	THR
2	H	170	LEU
2	H	171	MET
2	H	175	SER
2	H	184	LEU
2	H	190	HIS
2	H	191	ASN
2	H	192	THR
2	H	193	SER
2	H	194	LEU
2	H	196	VAL
2	H	203	GLU
2	H	214	THR
2	H	225	VAL
2	H	231	GLU
2	H	233	LEU
2	H	238	PHE
2	H	255	GLU
2	H	266	LEU
2	H	272	LEU
2	H	275	LEU
2	H	284	GLU
2	H	291	PHE
2	H	298	LEU
2	H	301	LEU
2	H	302	TYR
2	H	304	LEU
2	H	305	LEU

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Mol	Chain	Res	Type
2	H	313	LEU
2	H	315	GLN
2	H	316	LYS
2	H	325	THR
2	H	326	ARG
2	H	343	GLN
2	H	345	LYS
2	H	348	ARG
2	H	364	LEU
2	H	365	VAL
2	H	370	LEU
2	H	377	CYS
2	H	382	TYR
2	H	388	SER
2	H	389	ASP
2	H	390	ILE
2	H	394	SER
2	H	396	GLU
2	H	398	ILE
2	H	402	LEU
2	H	403	LYS
2	H	409	ARG
2	H	410	LEU
2	H	411	VAL
2	H	415	ARG
2	H	422	LEU
2	H	453	LEU
2	H	467	TRP
2	H	477	ASP
2	H	485	ARG
2	H	490	LEU
2	H	503	ARG
2	H	511	LYS
2	H	516	ARG
2	H	519	TRP
2	H	534	GLN
2	H	542	ILE
2	H	547	SER
2	H	548	ARG
2	H	568	ILE
2	H	579	THR
2	H	583	THR

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Mol	Chain	Res	Type
2	H	589	LYS
1	I	8	LEU
1	I	14	GLU
1	I	22	VAL
1	I	35	ASP
1	I	36	ASP
1	I	42	VAL
1	I	44	LEU
1	I	48	THR
1	I	81	ASP
1	I	125	ILE
1	I	150	GLU
1	I	153	ARG
1	I	160	GLU
2	J	16	THR
2	J	20	VAL
2	J	35	ASP
2	J	40	VAL
2	J	43	ARG
2	J	48	ASP
2	J	52	ARG
2	J	59	LEU
2	J	62	THR
2	J	64	THR
2	J	104	VAL
2	J	129	LEU
2	J	132	LEU
2	J	136	ARG
2	J	142	THR
2	J	159	ILE
2	J	166	ILE
2	J	168	THR
2	J	182	LYS
2	J	184	LEU
2	J	191	ASN
2	J	196	VAL
2	J	214	THR
2	J	217	ARG
2	J	220	ARG
2	J	225	VAL
2	J	227	VAL
2	J	230	PHE

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Mol	Chain	Res	Type
2	J	231	GLU
2	J	232	ILE
2	J	272	LEU
2	J	275	LEU
2	J	284	GLU
2	J	301	LEU
2	J	304	LEU
2	J	305	LEU
2	J	312	THR
2	J	314	ILE
2	J	327	ASN
2	J	341	CYS
2	J	349	ILE
2	J	360	ASP
2	J	367	GLN
2	J	377	CYS
2	J	390	ILE
2	J	400	THR
2	J	402	LEU
2	J	404	ASN
2	J	405	LEU
2	J	413	LEU
2	J	418	ARG
2	J	422	LEU
2	J	447	GLN
2	J	453	LEU
2	J	455	LEU
2	J	477	ASP
2	J	481	MET
2	J	490	LEU
2	J	495	MET
2	J	503	ARG
2	J	510	THR
2	J	516	ARG
2	J	519	TRP
2	J	528	THR
2	J	537	ARG
2	J	542	ILE
2	J	578	ARG
1	K	8	LEU
1	K	14	GLU
1	K	22	VAL

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Mol	Chain	Res	Type
1	K	29	ILE
1	K	35	ASP
1	K	36	ASP
1	K	42	VAL
1	K	44	LEU
1	K	48	THR
1	K	125	ILE
1	K	150	GLU
1	K	153	ARG
1	K	160	GLU
2	L	16	THR
2	L	20	VAL
2	L	35	ASP
2	L	40	VAL
2	L	43	ARG
2	L	48	ASP
2	L	52	ARG
2	L	59	LEU
2	L	62	THR
2	L	64	THR
2	L	104	VAL
2	L	123	ILE
2	L	129	LEU
2	L	132	LEU
2	L	136	ARG
2	L	142	THR
2	L	159	ILE
2	L	166	ILE
2	L	168	THR
2	L	182	LYS
2	L	184	LEU
2	L	191	ASN
2	L	193	SER
2	L	196	VAL
2	L	201	MET
2	L	214	THR
2	L	220	ARG
2	L	225	VAL
2	L	227	VAL
2	L	231	GLU
2	L	232	ILE
2	L	270	ARG

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Mol	Chain	Res	Type
2	L	272	LEU
2	L	275	LEU
2	L	284	GLU
2	L	301	LEU
2	L	304	LEU
2	L	305	LEU
2	L	312	THR
2	L	314	ILE
2	L	327	ASN
2	L	341	CYS
2	L	349	ILE
2	L	360	ASP
2	L	367	GLN
2	L	377	CYS
2	L	390	ILE
2	L	400	THR
2	L	402	LEU
2	L	404	ASN
2	L	405	LEU
2	L	413	LEU
2	L	418	ARG
2	L	422	LEU
2	L	447	GLN
2	L	453	LEU
2	L	455	LEU
2	L	477	ASP
2	L	481	MET
2	L	490	LEU
2	L	495	MET
2	L	503	ARG
2	L	510	THR
2	L	516	ARG
2	L	519	TRP
2	L	528	THR
2	L	537	ARG
2	L	542	ILE
2	L	578	ARG
1	M	8	LEU
1	M	14	GLU
1	M	22	VAL
1	M	35	ASP
1	M	36	ASP

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Mol	Chain	Res	Type
1	M	42	VAL
1	M	44	LEU
1	M	48	THR
1	M	81	ASP
1	M	125	ILE
1	M	150	GLU
1	M	153	ARG
1	M	160	GLU
2	N	16	THR
2	N	20	VAL
2	N	35	ASP
2	N	40	VAL
2	N	43	ARG
2	N	48	ASP
2	N	52	ARG
2	N	59	LEU
2	N	62	THR
2	N	64	THR
2	N	104	VAL
2	N	123	ILE
2	N	129	LEU
2	N	132	LEU
2	N	136	ARG
2	N	142	THR
2	N	159	ILE
2	N	166	ILE
2	N	168	THR
2	N	184	LEU
2	N	191	ASN
2	N	196	VAL
2	N	201	MET
2	N	214	THR
2	N	217	ARG
2	N	220	ARG
2	N	225	VAL
2	N	230	PHE
2	N	231	GLU
2	N	232	ILE
2	N	270	ARG
2	N	272	LEU
2	N	275	LEU
2	N	284	GLU

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Mol	Chain	Res	Type
2	N	301	LEU
2	N	304	LEU
2	N	305	LEU
2	N	312	THR
2	N	314	ILE
2	N	327	ASN
2	N	341	CYS
2	N	349	ILE
2	N	360	ASP
2	N	365	VAL
2	N	367	GLN
2	N	377	CYS
2	N	390	ILE
2	N	400	THR
2	N	402	LEU
2	N	404	ASN
2	N	405	LEU
2	N	418	ARG
2	N	422	LEU
2	N	447	GLN
2	N	453	LEU
2	N	455	LEU
2	N	477	ASP
2	N	481	MET
2	N	490	LEU
2	N	495	MET
2	N	503	ARG
2	N	510	THR
2	N	516	ARG
2	N	519	TRP
2	N	528	THR
2	N	537	ARG
2	N	542	ILE
2	N	578	ARG
1	O	8	LEU
1	O	14	GLU
1	O	22	VAL
1	O	29	ILE
1	O	35	ASP
1	O	36	ASP
1	O	42	VAL
1	O	44	LEU

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Mol	Chain	Res	Type
1	O	48	THR
1	O	81	ASP
1	O	125	ILE
1	O	150	GLU
1	O	153	ARG
1	O	160	GLU
2	P	16	THR
2	P	20	VAL
2	P	35	ASP
2	P	40	VAL
2	P	43	ARG
2	P	48	ASP
2	P	52	ARG
2	P	59	LEU
2	P	62	THR
2	P	64	THR
2	P	104	VAL
2	P	123	ILE
2	P	129	LEU
2	P	132	LEU
2	P	136	ARG
2	P	142	THR
2	P	159	ILE
2	P	168	THR
2	P	182	LYS
2	P	184	LEU
2	P	191	ASN
2	P	196	VAL
2	P	214	THR
2	P	220	ARG
2	P	225	VAL
2	P	230	PHE
2	P	231	GLU
2	P	232	ILE
2	P	270	ARG
2	P	272	LEU
2	P	275	LEU
2	P	284	GLU
2	P	301	LEU
2	P	304	LEU
2	P	305	LEU
2	P	312	THR

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Mol	Chain	Res	Type
2	P	314	ILE
2	P	327	ASN
2	P	341	CYS
2	P	349	ILE
2	P	360	ASP
2	P	367	GLN
2	P	377	CYS
2	P	390	ILE
2	P	400	THR
2	P	402	LEU
2	P	404	ASN
2	P	405	LEU
2	P	418	ARG
2	P	422	LEU
2	P	447	GLN
2	P	453	LEU
2	P	455	LEU
2	P	477	ASP
2	P	481	MET
2	P	490	LEU
2	P	495	MET
2	P	496	ARG
2	P	503	ARG
2	P	510	THR
2	P	516	ARG
2	P	519	TRP
2	P	528	THR
2	P	537	ARG
2	P	542	ILE
2	P	578	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (117) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	ASN
1	A	62	HIS
1	A	95	GLN
2	B	54	HIS
2	B	74	ASN
2	B	109	ASN
2	B	191	ASN
2	B	254	ASN

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Mol	Chain	Res	Type
2	B	294	GLN
2	B	319	ASN
2	B	327	ASN
2	B	343	GLN
2	B	460	GLN
2	B	489	ASN
2	B	564	HIS
2	B	567	HIS
1	C	46	ASN
1	C	62	HIS
2	D	54	HIS
2	D	74	ASN
2	D	109	ASN
2	D	190	HIS
2	D	191	ASN
2	D	254	ASN
2	D	294	GLN
2	D	319	ASN
2	D	327	ASN
2	D	343	GLN
2	D	460	GLN
2	D	489	ASN
2	D	564	HIS
2	D	567	HIS
1	E	46	ASN
1	E	62	HIS
1	E	95	GLN
2	F	54	HIS
2	F	74	ASN
2	F	109	ASN
2	F	113	GLN
2	F	191	ASN
2	F	254	ASN
2	F	327	ASN
2	F	343	GLN
2	F	460	GLN
2	F	489	ASN
2	F	564	HIS
2	F	567	HIS
1	G	46	ASN
1	G	62	HIS
2	H	23	GLN

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Mol	Chain	Res	Type
2	H	74	ASN
2	H	95	ASN
2	H	110	ASN
2	H	191	ASN
2	H	310	HIS
2	H	375	GLN
2	H	392	ASN
2	H	447	GLN
2	H	460	GLN
2	H	489	ASN
2	H	530	GLN
2	H	577	GLN
1	I	46	ASN
1	I	62	HIS
1	I	106	ASN
2	J	23	GLN
2	J	54	HIS
2	J	109	ASN
2	J	191	ASN
2	J	254	ASN
2	J	294	GLN
2	J	327	ASN
2	J	343	GLN
2	J	460	GLN
2	J	489	ASN
2	J	564	HIS
2	J	567	HIS
1	K	46	ASN
1	K	62	HIS
2	L	54	HIS
2	L	74	ASN
2	L	109	ASN
2	L	190	HIS
2	L	191	ASN
2	L	254	ASN
2	L	294	GLN
2	L	319	ASN
2	L	327	ASN
2	L	343	GLN
2	L	460	GLN
2	L	489	ASN
2	L	564	HIS

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Mol	Chain	Res	Type
2	L	567	HIS
1	M	46	ASN
1	M	62	HIS
2	N	54	HIS
2	N	109	ASN
2	N	191	ASN
2	N	254	ASN
2	N	327	ASN
2	N	343	GLN
2	N	460	GLN
2	N	489	ASN
2	N	564	HIS
2	N	567	HIS
1	O	46	ASN
1	O	62	HIS
2	P	54	HIS
2	P	109	ASN
2	P	191	ASN
2	P	254	ASN
2	P	327	ASN
2	P	343	GLN
2	P	460	GLN
2	P	489	ASN
2	P	564	HIS
2	P	567	HIS

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

40 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$
4	OGK	B	1100	-	25,25,25	7.08	10 (40%)	38,38,38	3.83	19 (50%)
5	PO4	B	1101	-	4,4,4	2.54	3 (75%)	6,6,6	0.32	0
5	PO4	B	1102	-	4,4,4	2.59	3 (75%)	6,6,6	0.31	0
5	PO4	B	1103	-	4,4,4	2.60	3 (75%)	6,6,6	0.31	0
5	PO4	B	1104	-	4,4,4	2.57	3 (75%)	6,6,6	0.31	0
5	PO4	D	1101	-	4,4,4	2.49	3 (75%)	6,6,6	0.35	0
5	PO4	D	1102	-	4,4,4	2.55	3 (75%)	6,6,6	0.30	0
5	PO4	D	1103	-	4,4,4	2.64	3 (75%)	6,6,6	0.32	0
5	PO4	D	1104	-	4,4,4	2.63	3 (75%)	6,6,6	0.32	0
4	OGK	D	2100	-	25,25,25	7.22	10 (40%)	38,38,38	3.80	18 (47%)
5	PO4	F	1101	-	4,4,4	2.56	3 (75%)	6,6,6	0.36	0
5	PO4	F	1102	-	4,4,4	2.76	3 (75%)	6,6,6	0.31	0
5	PO4	F	1103	-	4,4,4	2.63	3 (75%)	6,6,6	0.32	0
5	PO4	F	1104	-	4,4,4	2.74	3 (75%)	6,6,6	0.31	0
4	OGK	F	3100	-	25,25,25	7.09	11 (44%)	38,38,38	3.83	16 (42%)
5	PO4	H	1101	-	4,4,4	2.79	3 (75%)	6,6,6	0.31	0
5	PO4	H	1102	-	4,4,4	2.68	3 (75%)	6,6,6	0.32	0
5	PO4	H	1103	-	4,4,4	2.79	3 (75%)	6,6,6	0.31	0
5	PO4	H	1104	-	4,4,4	2.77	3 (75%)	6,6,6	0.35	0
4	OGK	H	4100	-	25,25,25	7.01	11 (44%)	38,38,38	3.64	19 (50%)
5	PO4	J	1101	-	4,4,4	2.60	3 (75%)	6,6,6	0.34	0
5	PO4	J	1102	-	4,4,4	2.66	3 (75%)	6,6,6	0.31	0
5	PO4	J	1103	-	4,4,4	2.50	3 (75%)	6,6,6	0.31	0
5	PO4	J	1104	-	4,4,4	2.52	3 (75%)	6,6,6	0.31	0
4	OGK	J	5100	-	25,25,25	7.22	11 (44%)	38,38,38	3.71	18 (47%)
5	PO4	L	1101	-	4,4,4	2.69	3 (75%)	6,6,6	0.33	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	PO4	L	1102	-	4,4,4	2.65	3 (75%)	6,6,6	0.30	0
5	PO4	L	1103	-	4,4,4	2.56	3 (75%)	6,6,6	0.32	0
5	PO4	L	1104	-	4,4,4	2.71	3 (75%)	6,6,6	0.31	0
4	OGK	L	6100	-	25,25,25	7.04	11 (44%)	38,38,38	3.68	17 (44%)
5	PO4	N	1101	-	4,4,4	2.90	3 (75%)	6,6,6	0.37	0
5	PO4	N	1102	-	4,4,4	2.68	3 (75%)	6,6,6	0.31	0
5	PO4	N	1103	-	4,4,4	2.57	3 (75%)	6,6,6	0.35	0
5	PO4	N	1104	-	4,4,4	2.71	3 (75%)	6,6,6	0.32	0
4	OGK	N	7100	-	25,25,25	6.81	10 (40%)	38,38,38	3.67	16 (42%)
5	PO4	P	1101	-	4,4,4	2.56	3 (75%)	6,6,6	0.33	0
5	PO4	P	1102	-	4,4,4	2.62	3 (75%)	6,6,6	0.31	0
5	PO4	P	1103	-	4,4,4	2.55	3 (75%)	6,6,6	0.32	0
5	PO4	P	1104	-	4,4,4	2.61	3 (75%)	6,6,6	0.32	0
4	OGK	P	8100	-	25,25,25	7.20	10 (40%)	38,38,38	3.72	18 (47%)

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	OGK	B	1100	-	-	0/19/52/52	0/0/3/3
5	PO4	B	1101	-	-	0/0/0/0	0/0/0/0
5	PO4	B	1102	-	-	0/0/0/0	0/0/0/0
5	PO4	B	1103	-	-	0/0/0/0	0/0/0/0
5	PO4	B	1104	-	-	0/0/0/0	0/0/0/0
5	PO4	D	1101	-	-	0/0/0/0	0/0/0/0
5	PO4	D	1102	-	-	0/0/0/0	0/0/0/0
5	PO4	D	1103	-	-	0/0/0/0	0/0/0/0
5	PO4	D	1104	-	-	0/0/0/0	0/0/0/0
4	OGK	D	2100	-	-	0/19/52/52	0/0/3/3
5	PO4	F	1101	-	-	0/0/0/0	0/0/0/0
5	PO4	F	1102	-	-	0/0/0/0	0/0/0/0
5	PO4	F	1103	-	-	0/0/0/0	0/0/0/0
5	PO4	F	1104	-	-	0/0/0/0	0/0/0/0
4	OGK	F	3100	-	-	0/19/52/52	0/0/3/3
5	PO4	H	1101	-	-	0/0/0/0	0/0/0/0
5	PO4	H	1102	-	-	0/0/0/0	0/0/0/0
5	PO4	H	1103	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PO4	H	1104	-	-	0/0/0/0	0/0/0/0
4	OGK	H	4100	-	1/1/9/10	0/19/52/52	0/0/3/3
5	PO4	J	1101	-	-	0/0/0/0	0/0/0/0
5	PO4	J	1102	-	-	0/0/0/0	0/0/0/0
5	PO4	J	1103	-	-	0/0/0/0	0/0/0/0
5	PO4	J	1104	-	-	0/0/0/0	0/0/0/0
4	OGK	J	5100	-	-	0/19/52/52	0/0/3/3
5	PO4	L	1101	-	-	0/0/0/0	0/0/0/0
5	PO4	L	1102	-	-	0/0/0/0	0/0/0/0
5	PO4	L	1103	-	-	0/0/0/0	0/0/0/0
5	PO4	L	1104	-	-	0/0/0/0	0/0/0/0
4	OGK	L	6100	-	-	0/19/52/52	0/0/3/3
5	PO4	N	1101	-	-	0/0/0/0	0/0/0/0
5	PO4	N	1102	-	-	0/0/0/0	0/0/0/0
5	PO4	N	1103	-	-	0/0/0/0	0/0/0/0
5	PO4	N	1104	-	-	0/0/0/0	0/0/0/0
4	OGK	N	7100	-	-	0/19/52/52	0/0/3/3
5	PO4	P	1101	-	-	0/0/0/0	0/0/0/0
5	PO4	P	1102	-	-	0/0/0/0	0/0/0/0
5	PO4	P	1103	-	-	0/0/0/0	0/0/0/0
5	PO4	P	1104	-	-	0/0/0/0	0/0/0/0
4	OGK	P	8100	-	-	0/19/52/52	0/0/3/3

All (180) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	J	5100	OGK	C13-C09	-23.21	1.27	1.51
4	D	2100	OGK	C13-C09	-23.11	1.27	1.51
4	P	8100	OGK	C13-C09	-22.84	1.28	1.51
4	F	3100	OGK	C13-C09	-22.78	1.28	1.51
4	B	1100	OGK	C13-C09	-22.53	1.28	1.51
4	L	6100	OGK	C13-C09	-22.51	1.28	1.51
4	H	4100	OGK	C13-C09	-22.28	1.28	1.51
4	N	7100	OGK	C13-C09	-21.47	1.29	1.51
4	P	8100	OGK	C09-C14	-18.55	1.33	1.52
4	D	2100	OGK	C09-C14	-18.36	1.33	1.52
4	J	5100	OGK	C09-C14	-18.13	1.34	1.52
4	F	3100	OGK	C09-C14	-18.04	1.34	1.52
4	B	1100	OGK	C09-C14	-17.94	1.34	1.52
4	L	6100	OGK	C09-C14	-17.80	1.34	1.52
4	H	4100	OGK	C09-C14	-17.77	1.34	1.52
4	J	5100	OGK	C09-C10	-17.09	1.13	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	N	7100	OGK	C09-C14	-17.01	1.35	1.52
4	D	2100	OGK	C09-C10	-16.98	1.13	1.52
4	P	8100	OGK	C09-C10	-16.95	1.14	1.52
4	L	6100	OGK	C09-C10	-16.65	1.14	1.52
4	B	1100	OGK	C09-C10	-16.44	1.15	1.52
4	F	3100	OGK	C09-C10	-16.40	1.15	1.52
4	H	4100	OGK	C09-C10	-16.15	1.15	1.52
4	N	7100	OGK	C09-C10	-16.10	1.15	1.52
4	B	1100	OGK	C06-N08	7.05	1.49	1.34
4	F	3100	OGK	C06-N08	6.92	1.49	1.34
4	P	8100	OGK	C06-N08	6.87	1.49	1.34
4	B	1100	OGK	C13-C14	-6.84	1.31	1.50
4	D	2100	OGK	C13-C14	-6.70	1.32	1.50
4	N	7100	OGK	C06-N08	6.64	1.48	1.34
4	D	2100	OGK	C06-N08	6.60	1.48	1.34
4	H	4100	OGK	C13-C14	-6.59	1.32	1.50
4	J	5100	OGK	C13-C14	-6.54	1.32	1.50
4	L	6100	OGK	C13-C14	-6.49	1.32	1.50
4	F	3100	OGK	C13-C14	-6.42	1.33	1.50
4	P	8100	OGK	C13-C14	-6.33	1.33	1.50
4	J	5100	OGK	C06-N08	6.26	1.47	1.34
4	L	6100	OGK	C06-N08	6.22	1.47	1.34
4	H	4100	OGK	C06-N08	6.19	1.47	1.34
4	N	7100	OGK	C13-C14	-6.08	1.33	1.50
4	H	4100	OGK	C23-C22	4.81	1.61	1.53
4	N	7100	OGK	C23-C22	4.22	1.60	1.53
4	N	7100	OGK	O12-C10	4.01	1.45	1.30
5	N	1101	PO4	P-O3	3.92	1.68	1.52
4	H	4100	OGK	O12-C10	3.86	1.44	1.30
4	B	1100	OGK	O12-C10	3.60	1.43	1.30
5	H	1104	PO4	P-O3	3.52	1.67	1.52
4	F	3100	OGK	O12-C10	3.52	1.43	1.30
4	P	8100	OGK	C23-C22	3.51	1.59	1.53
5	N	1101	PO4	P-O4	3.46	1.66	1.52
5	N	1104	PO4	P-O4	3.46	1.66	1.52
4	L	6100	OGK	O12-C10	3.42	1.42	1.30
4	F	3100	OGK	C18-C17	-3.42	1.44	1.54
5	H	1103	PO4	P-O3	3.40	1.66	1.52
5	H	1101	PO4	P-O3	3.38	1.66	1.52
4	L	6100	OGK	C23-C22	3.38	1.59	1.53
4	P	8100	OGK	O12-C10	3.37	1.42	1.30
5	L	1104	PO4	P-O4	3.37	1.66	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1100	OGK	C18-C17	-3.36	1.44	1.54
4	D	2100	OGK	O12-C10	3.35	1.42	1.30
4	J	5100	OGK	O12-C10	3.35	1.42	1.30
5	B	1103	PO4	P-O3	3.34	1.66	1.52
5	H	1104	PO4	P-O4	3.32	1.66	1.52
5	H	1101	PO4	P-O4	3.32	1.66	1.52
5	H	1102	PO4	P-O4	3.31	1.66	1.52
5	F	1103	PO4	P-O4	3.29	1.66	1.52
4	P	8100	OGK	C18-C17	-3.28	1.45	1.54
5	L	1101	PO4	P-O3	3.27	1.66	1.52
5	N	1102	PO4	P-O4	3.27	1.66	1.52
5	F	1102	PO4	P-O3	3.26	1.66	1.52
4	B	1100	OGK	C23-C22	3.26	1.58	1.53
5	P	1102	PO4	P-O3	3.26	1.65	1.52
5	J	1102	PO4	P-O3	3.25	1.65	1.52
5	F	1104	PO4	P-O4	3.24	1.65	1.52
4	J	5100	OGK	C18-C17	-3.24	1.45	1.54
5	F	1104	PO4	P-O3	3.24	1.65	1.52
5	D	1104	PO4	P-O3	3.23	1.65	1.52
5	J	1101	PO4	P-O4	3.23	1.65	1.52
5	L	1101	PO4	P-O4	3.22	1.65	1.52
5	H	1103	PO4	P-O4	3.22	1.65	1.52
5	B	1101	PO4	P-O3	3.20	1.65	1.52
5	L	1102	PO4	P-O3	3.20	1.65	1.52
5	N	1103	PO4	P-O3	3.19	1.65	1.52
5	N	1104	PO4	P-O3	3.19	1.65	1.52
4	D	2100	OGK	C23-C22	3.17	1.58	1.53
4	J	5100	OGK	C23-C22	3.15	1.58	1.53
5	N	1102	PO4	P-O3	3.15	1.65	1.52
4	D	2100	OGK	C18-C17	-3.14	1.45	1.54
5	P	1104	PO4	P-O4	3.14	1.65	1.52
5	J	1101	PO4	P-O3	3.14	1.65	1.52
5	L	1103	PO4	P-O3	3.14	1.65	1.52
4	H	4100	OGK	C18-C17	-3.12	1.45	1.54
5	F	1102	PO4	P-O4	3.12	1.65	1.52
5	L	1104	PO4	P-O3	3.11	1.65	1.52
4	N	7100	OGK	C18-C17	-3.11	1.45	1.54
5	F	1103	PO4	P-O3	3.11	1.65	1.52
5	D	1102	PO4	P-O4	3.11	1.65	1.52
5	B	1102	PO4	P-O4	3.10	1.65	1.52
5	D	1104	PO4	P-O4	3.10	1.65	1.52
5	P	1104	PO4	P-O3	3.07	1.65	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	J	1102	PO4	P-O4	3.06	1.65	1.52
5	P	1103	PO4	P-O4	3.06	1.65	1.52
5	P	1101	PO4	P-O3	3.06	1.65	1.52
5	P	1103	PO4	P-O3	3.05	1.65	1.52
5	J	1104	PO4	P-O3	3.05	1.65	1.52
5	H	1102	PO4	P-O3	3.04	1.65	1.52
5	B	1104	PO4	P-O3	3.03	1.65	1.52
5	P	1102	PO4	P-O4	3.03	1.65	1.52
5	N	1103	PO4	P-O4	3.03	1.65	1.52
5	D	1101	PO4	P-O4	3.03	1.65	1.52
5	L	1102	PO4	P-O4	3.02	1.65	1.52
5	D	1103	PO4	P-O3	3.01	1.65	1.52
5	B	1102	PO4	P-O3	2.97	1.64	1.52
5	J	1103	PO4	P-O3	2.97	1.64	1.52
5	F	1101	PO4	P-O3	2.96	1.64	1.52
5	D	1103	PO4	P-O4	2.96	1.64	1.52
5	B	1104	PO4	P-O4	2.94	1.64	1.52
5	P	1101	PO4	P-O4	2.94	1.64	1.52
5	F	1101	PO4	P-O4	2.93	1.64	1.52
4	L	6100	OGK	C18-C17	-2.90	1.46	1.54
5	D	1101	PO4	P-O3	2.88	1.64	1.52
4	B	1100	OGK	C09-N08	2.86	1.50	1.45
4	N	7100	OGK	C09-N08	2.85	1.50	1.45
5	J	1104	PO4	P-O4	2.81	1.64	1.52
4	F	3100	OGK	C23-C22	2.76	1.58	1.53
5	L	1103	PO4	P-O4	2.75	1.63	1.52
4	L	6100	OGK	C09-N08	2.75	1.50	1.45
5	J	1103	PO4	P-O4	2.75	1.63	1.52
5	D	1102	PO4	P-O3	2.73	1.63	1.52
5	B	1103	PO4	P-O4	2.73	1.63	1.52
5	F	1102	PO4	P-O1	2.71	1.66	1.52
4	J	5100	OGK	C05-C06	2.67	1.56	1.51
4	D	2100	OGK	C05-C06	2.67	1.56	1.51
4	F	3100	OGK	C09-N08	2.67	1.50	1.45
5	B	1101	PO4	P-O4	2.66	1.63	1.52
4	J	5100	OGK	C09-N08	2.63	1.50	1.45
5	F	1104	PO4	P-O1	2.63	1.66	1.52
5	J	1104	PO4	P-O1	2.63	1.66	1.52
4	L	6100	OGK	C05-C06	2.58	1.56	1.51
5	L	1104	PO4	P-O1	2.57	1.66	1.52
5	H	1103	PO4	P-O1	2.54	1.66	1.52
4	J	5100	OGK	O07-C06	-2.52	1.18	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	4100	OGK	C05-C06	2.50	1.56	1.51
5	B	1102	PO4	P-O1	2.49	1.65	1.52
5	N	1102	PO4	P-O1	2.48	1.65	1.52
5	N	1103	PO4	P-O1	2.48	1.65	1.52
5	J	1103	PO4	P-O1	2.47	1.65	1.52
4	D	2100	OGK	O07-C06	-2.47	1.18	1.23
5	L	1103	PO4	P-O1	2.47	1.65	1.52
5	D	1103	PO4	P-O1	2.46	1.65	1.52
5	P	1104	PO4	P-O1	2.46	1.65	1.52
5	D	1102	PO4	P-O1	2.44	1.65	1.52
5	N	1101	PO4	P-O1	2.43	1.65	1.52
5	H	1102	PO4	P-O1	2.43	1.65	1.52
5	N	1104	PO4	P-O1	2.43	1.65	1.52
4	H	4100	OGK	C09-N08	2.43	1.50	1.45
5	P	1103	PO4	P-O1	2.42	1.65	1.52
5	J	1102	PO4	P-O1	2.42	1.65	1.52
5	F	1101	PO4	P-O1	2.42	1.65	1.52
5	F	1103	PO4	P-O1	2.42	1.65	1.52
5	D	1104	PO4	P-O1	2.41	1.65	1.52
5	B	1104	PO4	P-O1	2.40	1.65	1.52
5	H	1104	PO4	P-O1	2.40	1.65	1.52
5	H	1101	PO4	P-O1	2.40	1.65	1.52
4	H	4100	OGK	O07-C06	-2.40	1.18	1.23
5	P	1102	PO4	P-O1	2.40	1.65	1.52
5	B	1103	PO4	P-O1	2.40	1.65	1.52
5	P	1101	PO4	P-O1	2.39	1.65	1.52
4	P	8100	OGK	C05-C06	2.38	1.56	1.51
5	L	1102	PO4	P-O1	2.38	1.65	1.52
4	L	6100	OGK	O07-C06	-2.33	1.18	1.23
5	B	1101	PO4	P-O1	2.32	1.64	1.52
5	J	1101	PO4	P-O1	2.28	1.64	1.52
5	L	1101	PO4	P-O1	2.27	1.64	1.52
4	B	1100	OGK	O07-C06	-2.23	1.19	1.23
5	D	1101	PO4	P-O1	2.18	1.64	1.52
4	F	3100	OGK	C05-C06	2.13	1.55	1.51
4	P	8100	OGK	C09-N08	2.09	1.49	1.45
4	N	7100	OGK	O07-C06	-2.01	1.19	1.23
4	F	3100	OGK	O07-C06	-2.01	1.19	1.23

All (141) bond angle outliers are listed below:

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1100	OGK	C15-C14-C09	-16.31	99.15	122.07
4	F	3100	OGK	C15-C14-C09	-16.25	99.23	122.07
4	D	2100	OGK	C15-C14-C09	-15.66	100.06	122.07
4	J	5100	OGK	C15-C14-C09	-15.40	100.43	122.07
4	L	6100	OGK	C15-C14-C09	-15.27	100.61	122.07
4	P	8100	OGK	C15-C14-C09	-15.18	100.73	122.07
4	N	7100	OGK	C15-C14-C09	-14.97	101.02	122.07
4	H	4100	OGK	C15-C14-C09	-14.36	101.89	122.07
4	B	1100	OGK	C18-C17-C05	-8.13	100.00	118.31
4	F	3100	OGK	C18-C17-C05	-7.92	100.47	118.31
4	P	8100	OGK	C18-C17-C05	-7.78	100.80	118.31
4	N	7100	OGK	C18-C17-C05	-7.60	101.19	118.31
4	D	2100	OGK	C18-C17-C05	-7.60	101.21	118.31
4	L	6100	OGK	C18-C17-C05	-7.16	102.19	118.31
4	J	5100	OGK	C18-C17-C05	-7.04	102.47	118.31
4	J	5100	OGK	C13-C14-C09	-6.83	57.22	60.40
4	F	3100	OGK	C13-C14-C09	-6.59	57.33	60.40
4	P	8100	OGK	C13-C14-C09	-6.50	57.37	60.40
4	D	2100	OGK	C13-C14-C09	-6.37	57.43	60.40
4	L	6100	OGK	C13-C14-C09	-6.33	57.45	60.40
4	N	7100	OGK	C13-C14-C09	-6.21	57.51	60.40
4	H	4100	OGK	C13-C14-C09	-5.95	57.63	60.40
4	B	1100	OGK	C13-C14-C09	-5.79	57.70	60.40
4	H	4100	OGK	C18-C17-C05	-5.68	105.53	118.31
4	N	7100	OGK	C04-C05-C17	5.55	121.81	109.95
4	L	6100	OGK	C05-C04-C03	5.21	119.91	109.75
4	B	1100	OGK	C04-C05-C17	5.18	121.01	109.95
4	P	8100	OGK	C04-C05-C17	5.15	120.95	109.95
4	J	5100	OGK	C04-C05-C17	5.09	120.82	109.95
4	H	4100	OGK	C23-C03-C02	5.08	121.43	112.79
4	P	8100	OGK	C05-C04-C03	5.06	119.63	109.75
4	F	3100	OGK	C04-C05-C17	4.99	120.60	109.95
4	L	6100	OGK	C04-C05-C17	4.97	120.58	109.95
4	D	2100	OGK	C05-C04-C03	4.96	119.42	109.75
4	D	2100	OGK	C04-C05-C17	4.87	120.35	109.95
4	J	5100	OGK	C05-C04-C03	4.66	118.84	109.75
4	B	1100	OGK	C05-C04-C03	4.61	118.74	109.75
4	H	4100	OGK	C14-C09-C10	4.58	128.15	117.31
4	H	4100	OGK	C17-C05-C06	4.54	119.81	111.64
4	F	3100	OGK	C05-C04-C03	4.51	118.55	109.75
4	D	2100	OGK	C14-C09-C10	4.49	127.92	117.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	N	7100	OGK	C18-C17-C22	4.42	109.22	103.97
4	H	4100	OGK	C13-C09-N08	-4.35	112.35	117.67
4	B	1100	OGK	C14-C09-C10	4.29	127.46	117.31
4	N	7100	OGK	C14-C09-C10	4.28	127.44	117.31
4	N	7100	OGK	C05-C04-C03	4.27	118.08	109.75
4	F	3100	OGK	C17-C05-C06	4.24	119.27	111.64
4	H	4100	OGK	C04-C05-C06	4.19	120.30	110.98
4	B	1100	OGK	O07-C06-C05	-4.16	116.69	121.51
4	F	3100	OGK	C14-C09-C10	4.10	127.00	117.31
4	P	8100	OGK	C13-C09-C14	4.04	61.17	59.11
4	P	8100	OGK	C14-C09-C10	4.00	126.76	117.31
4	L	6100	OGK	C14-C09-C10	3.96	126.66	117.31
4	P	8100	OGK	C17-C05-C06	3.92	118.69	111.64
4	L	6100	OGK	C17-C05-C06	3.86	118.58	111.64
4	F	3100	OGK	C18-C17-C22	3.78	108.46	103.97
4	J	5100	OGK	C14-C09-C10	3.78	126.24	117.31
4	B	1100	OGK	C17-C05-C06	3.74	118.38	111.64
4	J	5100	OGK	C18-C17-C22	3.74	108.41	103.97
4	D	2100	OGK	C13-C09-N08	-3.71	113.14	117.67
4	D	2100	OGK	C17-C05-C06	3.70	118.31	111.64
4	J	5100	OGK	C17-C05-C06	3.66	118.22	111.64
4	D	2100	OGK	C04-C03-C02	-3.65	106.58	112.79
4	L	6100	OGK	C04-C05-C06	3.59	118.97	110.98
4	H	4100	OGK	C04-C05-C17	3.54	117.52	109.95
4	D	2100	OGK	C04-C05-C06	3.49	118.74	110.98
4	F	3100	OGK	C13-C09-C14	3.48	60.89	59.11
4	H	4100	OGK	C22-C23-C03	3.41	116.40	109.75
4	J	5100	OGK	C13-C09-C14	3.39	60.84	59.11
4	P	8100	OGK	C04-C05-C06	3.38	118.50	110.98
4	F	3100	OGK	C13-C09-N08	-3.37	113.54	117.67
4	H	4100	OGK	C14-C09-N08	-3.36	108.61	117.92
4	H	4100	OGK	O12-C10-C09	3.32	120.18	113.25
4	J	5100	OGK	C04-C05-C06	3.30	118.33	110.98
4	J	5100	OGK	C04-C03-C02	-3.29	107.20	112.79
4	D	2100	OGK	C18-C17-C22	3.29	107.87	103.97
4	L	6100	OGK	C04-C03-C02	-3.27	107.22	112.79
4	H	4100	OGK	C04-C03-C23	3.25	115.98	109.72
4	N	7100	OGK	C13-C09-C14	3.23	60.76	59.11
4	N	7100	OGK	C17-C05-C06	3.17	117.35	111.64
4	P	8100	OGK	C13-C09-N08	-3.17	113.79	117.67
4	L	6100	OGK	C13-C09-N08	-3.17	113.80	117.67
4	F	3100	OGK	C14-C09-N08	-3.16	109.17	117.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	2100	OGK	C14-C09-N08	-3.15	109.19	117.92
4	B	1100	OGK	C04-C05-C06	3.13	117.95	110.98
4	N	7100	OGK	O11-C10-C09	-3.13	118.42	122.85
4	H	4100	OGK	C18-C17-C22	3.12	107.68	103.97
4	N	7100	OGK	C04-C05-C06	3.06	117.80	110.98
4	F	3100	OGK	C04-C05-C06	3.05	117.76	110.98
4	L	6100	OGK	O12-C10-C09	3.03	119.58	113.25
4	N	7100	OGK	C13-C09-N08	-3.03	113.97	117.67
4	D	2100	OGK	C13-C09-C14	3.02	60.65	59.11
4	H	4100	OGK	O11-C10-C09	-3.02	118.57	122.85
4	L	6100	OGK	C13-C09-C14	3.02	60.65	59.11
4	P	8100	OGK	C04-C03-C02	-3.01	107.67	112.79
4	D	2100	OGK	O12-C10-C09	3.00	119.53	113.25
4	P	8100	OGK	O12-C10-C09	3.00	119.51	113.25
4	P	8100	OGK	C18-C17-C22	2.98	107.50	103.97
4	B	1100	OGK	C13-C09-N08	-2.97	114.03	117.67
4	P	8100	OGK	C14-C09-N08	-2.97	109.70	117.92
4	D	2100	OGK	O11-C10-C09	-2.96	118.65	122.85
4	F	3100	OGK	O12-C10-C09	2.96	119.43	113.25
4	J	5100	OGK	C13-C09-N08	-2.95	114.07	117.67
4	J	5100	OGK	C14-C09-N08	-2.91	109.85	117.92
4	J	5100	OGK	O12-C10-C09	2.91	119.34	113.25
4	L	6100	OGK	C14-C09-N08	-2.90	109.88	117.92
4	B	1100	OGK	C04-C03-C02	-2.90	107.86	112.79
4	N	7100	OGK	C14-C09-N08	-2.89	109.93	117.92
4	B	1100	OGK	C14-C09-N08	-2.86	110.00	117.92
4	B	1100	OGK	C05-C06-N08	2.82	120.19	116.28
4	J	5100	OGK	O11-C10-C09	-2.79	118.90	122.85
4	L	6100	OGK	C18-C17-C22	2.79	107.28	103.97
4	B	1100	OGK	O12-C10-C09	2.77	119.03	113.25
4	B	1100	OGK	C15-C14-C13	-2.67	117.56	121.45
4	N	7100	OGK	C16-C15-C14	2.62	117.74	112.94
4	H	4100	OGK	C13-C09-C14	2.61	60.44	59.11
4	H	4100	OGK	C05-C04-C03	2.60	114.82	109.75
4	P	8100	OGK	O07-C06-C05	-2.59	118.50	121.51
4	J	5100	OGK	C05-C06-N08	2.55	119.82	116.28
4	N	7100	OGK	O12-C10-C09	2.55	118.57	113.25
4	B	1100	OGK	C09-C13-C14	2.51	62.10	60.84
4	F	3100	OGK	C04-C03-C02	-2.51	108.52	112.79
4	P	8100	OGK	O11-C10-C09	-2.50	119.31	122.85
4	D	2100	OGK	C05-C06-N08	2.45	119.69	116.28
4	B	1100	OGK	O11-C10-C09	-2.44	119.39	122.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	3100	OGK	O07-C06-C05	-2.40	118.73	121.51
4	D	2100	OGK	C15-C14-C13	-2.32	118.07	121.45
4	L	6100	OGK	O11-C10-C09	-2.29	119.61	122.85
4	L	6100	OGK	C05-C06-N08	2.28	119.44	116.28
4	J	5100	OGK	C09-C13-C14	2.20	61.94	60.84
4	B	1100	OGK	C18-C17-C22	2.19	106.57	103.97
4	H	4100	OGK	C09-C13-C14	2.18	61.93	60.84
4	F	3100	OGK	O11-C10-C09	-2.17	119.77	122.85
4	N	7100	OGK	C04-C03-C02	-2.17	109.10	112.79
4	J	5100	OGK	O07-C06-N08	-2.17	117.86	123.00
4	P	8100	OGK	C16-C15-C14	2.16	116.89	112.94
4	D	2100	OGK	C09-C13-C14	2.15	61.92	60.84
4	B	1100	OGK	C13-C09-C14	2.14	60.20	59.11
4	H	4100	OGK	C23-C22-C17	2.12	119.89	112.92
4	L	6100	OGK	C09-C13-C14	2.11	61.90	60.84
4	P	8100	OGK	C05-C06-N08	2.04	119.12	116.28

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	H	4100	OGK	C03

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	144/160 (90%)	0.26	1 (0%) 84 43	50, 100, 151, 168	0
1	C	144/160 (90%)	0.22	0 100 100	50, 102, 153, 170	0
1	E	144/160 (90%)	0.15	0 100 100	54, 101, 151, 171	0
1	G	144/160 (90%)	0.02	0 100 100	49, 96, 149, 170	0
1	I	144/160 (90%)	0.12	0 100 100	50, 100, 150, 170	0
1	K	144/160 (90%)	0.23	1 (0%) 84 43	53, 102, 152, 170	0
1	M	144/160 (90%)	0.34	2 (1%) 72 26	55, 102, 152, 175	0
1	O	144/160 (90%)	0.28	2 (1%) 72 26	51, 101, 151, 173	0
2	B	568/592 (95%)	-0.13	4 (0%) 84 43	38, 76, 136, 201	0
2	D	568/592 (95%)	-0.08	2 (0%) 90 58	38, 75, 135, 200	0
2	F	568/592 (95%)	-0.09	10 (1%) 65 22	40, 79, 136, 202	0
2	H	562/592 (94%)	-0.18	3 (0%) 88 52	34, 73, 127, 191	0
2	J	568/592 (95%)	-0.12	6 (1%) 77 32	39, 77, 136, 204	0
2	L	568/592 (95%)	-0.13	4 (0%) 84 43	41, 79, 135, 202	0
2	N	568/592 (95%)	0.18	12 (2%) 60 19	41, 82, 137, 206	0
2	P	568/592 (95%)	-0.09	7 (1%) 75 30	40, 80, 138, 204	0
3	Q	18/21 (85%)	-0.10	0 100 100	72, 93, 116, 143	0
3	R	18/21 (85%)	-0.08	0 100 100	71, 88, 115, 145	0
3	S	18/21 (85%)	-0.24	0 100 100	73, 93, 117, 143	0
3	U	18/21 (85%)	-0.17	0 100 100	71, 88, 112, 143	0
3	V	18/21 (85%)	-0.03	1 (5%) 24 6	75, 93, 115, 143	0
3	W	18/21 (85%)	0.17	1 (5%) 24 6	80, 96, 119, 143	0
3	X	18/21 (85%)	-0.11	1 (5%) 24 6	76, 91, 119, 143	0
All	All	5816/6163 (94%)	-0.02	57 (0%) 79 34	34, 82, 141, 206	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	N	591	PRO	6.0
2	N	590	GLU	5.9
2	N	592	ILE	4.6
2	B	549	ARG	4.4
2	P	362	GLU	4.1
2	P	592	ILE	4.0
2	F	355	GLU	3.5
2	N	355	GLU	3.5
2	J	360	ASP	3.3
2	F	356	GLN	3.2
2	N	356	GLN	3.0
2	L	359	GLU	3.0
2	J	592	ILE	3.0
3	V	200	GLU	2.9
2	F	592	ILE	2.9
2	L	592	ILE	2.9
2	N	359	GLU	2.8
1	A	24	LEU	2.7
3	W	200	GLU	2.7
2	H	355	GLU	2.7
2	P	591	PRO	2.7
2	D	360	ASP	2.6
2	P	548	ARG	2.6
2	J	549	ARG	2.6
2	H	356	GLN	2.6
2	N	266	LEU	2.5
2	J	358	MET	2.5
2	B	592	ILE	2.5
1	M	80	ASP	2.5
2	H	362	GLU	2.4
2	F	590	GLU	2.4
2	B	356	GLN	2.4
2	J	548	ARG	2.4
2	P	356	GLN	2.4
2	N	358	MET	2.3
2	F	416	GLU	2.3
2	J	361	GLU	2.3
2	N	360	ASP	2.3
3	X	200	GLU	2.3
2	B	355	GLU	2.2
2	F	358	MET	2.2
2	P	547	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	O	5	LYS	2.2
1	K	6	ILE	2.2
2	N	527	MET	2.2
1	O	80	ASP	2.2
2	F	591	PRO	2.2
2	N	548	ARG	2.1
2	P	590	GLU	2.1
2	F	546	PRO	2.1
1	M	115	ASP	2.1
2	L	356	GLN	2.1
2	N	373	LEU	2.1
2	D	592	ILE	2.1
2	L	360	ASP	2.0
2	F	359	GLU	2.0
2	F	547	SER	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	PO4	J	1102	5/5	0.39	17.65	67,86,95,104	0
5	PO4	F	1102	5/5	0.33	7.86	64,90,96,109	0
5	PO4	L	1102	5/5	0.38	5.81	73,82,96,100	0
5	PO4	H	1104	5/5	0.29	5.41	54,57,114,134	0
5	PO4	H	1103	5/5	0.18	5.20	38,47,52,77	0
5	PO4	N	1102	5/5	0.34	4.56	78,90,92,113	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	PO4	L	1103	5/5	0.32	3.99	52,56,69,89	0
5	PO4	L	1101	5/5	0.32	3.48	48,58,74,76	0
5	PO4	H	1101	5/5	0.20	3.17	36,45,63,88	0
5	PO4	B	1101	5/5	0.24	3.16	58,59,65,75	0
5	PO4	H	1102	5/5	0.29	3.08	58,59,72,88	0
5	PO4	J	1101	5/5	0.25	3.00	50,55,74,79	0
5	PO4	B	1102	5/5	0.26	2.71	70,84,93,100	0
5	PO4	F	1103	5/5	0.30	2.48	53,57,73,86	0
5	PO4	P	1101	5/5	0.24	2.40	57,62,74,75	0
5	PO4	B	1104	5/5	0.25	2.28	68,79,102,119	0
5	PO4	N	1103	5/5	0.33	2.21	56,56,74,85	0
5	PO4	F	1101	5/5	0.23	2.17	52,58,68,88	0
5	PO4	B	1103	5/5	0.28	2.15	54,54,61,76	0
5	PO4	D	1101	5/5	0.28	1.76	51,55,63,80	0
5	PO4	J	1104	5/5	0.26	1.69	76,81,112,117	0
5	PO4	L	1104	5/5	0.24	1.64	76,78,119,123	0
5	PO4	D	1103	5/5	0.30	1.63	48,52,55,81	0
5	PO4	D	1102	5/5	0.23	1.61	66,79,93,93	0
4	OGK	H	4100	23/23	0.26	1.38	44,71,85,98	0
5	PO4	J	1103	5/5	0.24	1.15	53,53,64,77	0
4	OGK	P	8100	23/23	0.26	0.95	64,75,88,90	0
5	PO4	P	1104	5/5	0.21	0.90	73,75,110,121	0
5	PO4	D	1104	5/5	0.23	0.81	72,73,107,122	0
4	OGK	J	5100	23/23	0.25	0.65	59,67,88,93	0
5	PO4	N	1104	5/5	0.24	0.63	82,83,118,122	0
5	PO4	P	1103	5/5	0.22	0.58	50,58,73,85	0
4	OGK	D	2100	23/23	0.26	0.54	60,72,81,86	0
4	OGK	B	1100	23/23	0.23	0.46	62,73,87,93	0
4	OGK	F	3100	23/23	0.23	0.46	64,72,91,94	0
5	PO4	N	1101	5/5	0.29	0.39	51,59,80,82	0
4	OGK	L	6100	23/23	0.22	0.29	62,71,90,92	0
4	OGK	N	7100	23/23	0.29	0.23	65,79,100,104	0
5	PO4	F	1104	5/5	0.21	0.14	72,81,117,120	0
5	PO4	P	1102	5/5	0.16	-2.18	72,87,98,108	0

6.5 Other polymers ⓘ

There are no such residues in this entry.