



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 12:04 AM GMT

PDB ID : 1VF5  
Title : Crystal Structure of Cytochrome b6f Complex from M.laminosus  
Authors : Kurisu, G.; Zhang, H.; Smith, J.L.; Cramer, W.A.  
Deposited on : 2004-04-08  
Resolution : 3.00 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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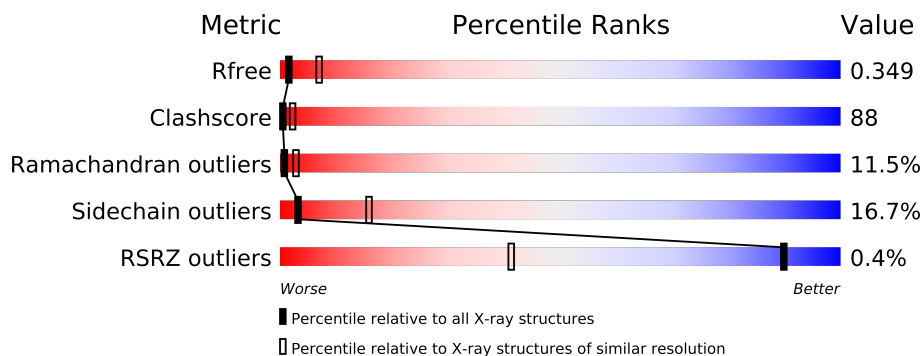
The following versions of software and data (see [references](#)) were used in the production of this report:

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

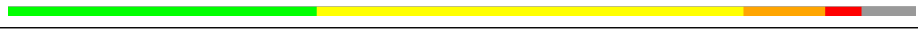
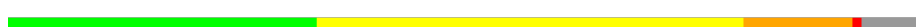
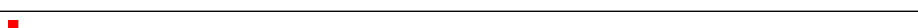

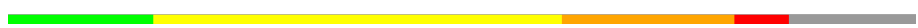
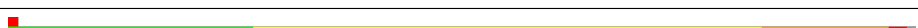

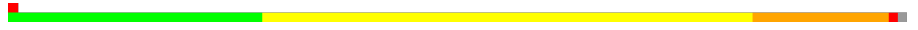
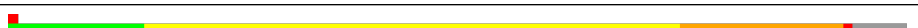


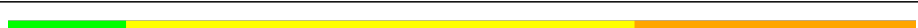
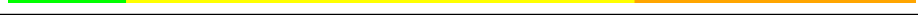
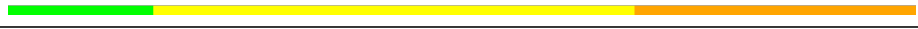
The reported resolution of this entry is 3.00 Å.

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



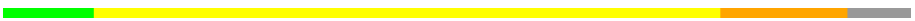
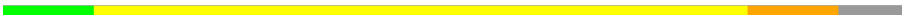
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	66092	1216 (3.00-3.00)
Clashscore	79885	1594 (3.00-3.00)
Ramachandran outliers	78287	1537 (3.00-3.00)
Sidechain outliers	78261	1540 (3.00-3.00)
RSRZ outliers	66119	1217 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	215	
1	N	215	
2	B	160	
2	O	160	
3	C	289	
3	P	289	
4	D	179	
4	Q	179	
5	E	32	
5	R	32	
6	F	35	
6	S	35	
7	G	37	
7	T	37	

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Mol	Chain	Length	Quality of chain
8	H	29	
8	U	29	

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

Mol	Type	Chain	Res	Geometry	Electron density
10	TDS	N	1304	-	X
11	PL9	A	305	-	X
13	CLA	B	201	-	X
15	BCR	E	101	-	X
15	BCR	R	1101	-	X
9	HEM	A	301	-	X

## 2 Entry composition

There are 16 unique types of molecules in this entry. The entry contains 15091 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 CYTOCHROME B6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	202	Total	C	N	O	S	0	0	0
			1593	1062	253	268	10			
1	N	202	Total	C	N	O	S	0	0	0
			1593	1062	253	268	10			

- Molecule 2 is a protein called SUBUNIT IV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	138	Total	C	N	O	S	0	0	0
			1075	727	165	178	5			
2	O	138	Total	C	N	O	S	0	0	0
			1075	727	165	178	5			

- Molecule 3 is a protein called CYTOCHROME F.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	286	Total	C	N	O	S	0	0	0
			2200	1406	366	421	7			
3	P	286	Total	C	N	O	S	0	0	0
			2200	1406	366	421	7			

- Molecule 4 is a protein called RIESKE IRON-SULFUR PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	168	Total	C	N	O	S	0	0	0
			1280	815	223	235	7			
4	Q	168	Total	C	N	O	S	0	0	0
			1280	815	223	235	7			

- Molecule 5 is a protein called PROTEIN PET L.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	32	Total	C	N	O	S	0	0	0
			248	179	34	34	1			
5	R	32	Total	C	N	O	S	0	0	0
			248	179	34	34	1			

- Molecule 6 is a protein called PROTEIN PET M.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	33	Total	C	N	O	S	0	0	0
			251	170	36	43	2			
6	S	35	Total	C	N	O	S	0	0	0
			270	181	39	48	2			

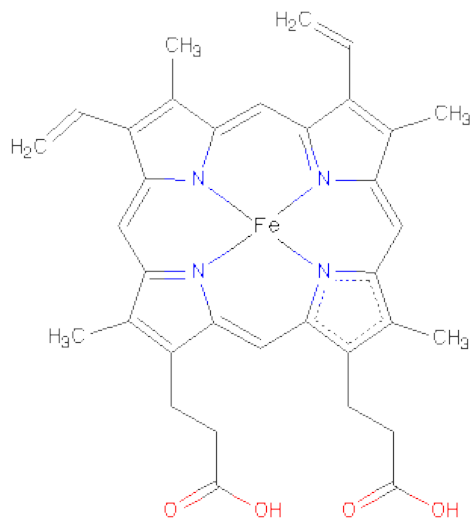
- Molecule 7 is a protein called PROTEIN PET G.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	G	23	Total	C	N	O	0	0	0
			184	126	29	29			
7	T	27	Total	C	N	O	0	0	0
			216	146	34	36			

- Molecule 8 is a protein called PROTEIN PET N.

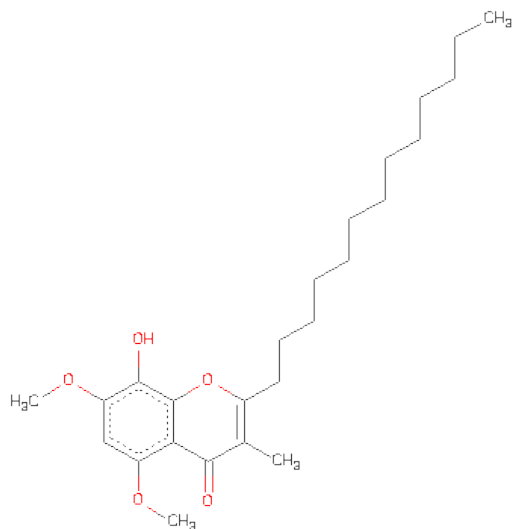
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	27	Total	C	N	O	S	0	0	0
			214	146	34	33	1			
8	U	27	Total	C	N	O	S	0	0	0
			214	146	34	33	1			

- Molecule 9 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).



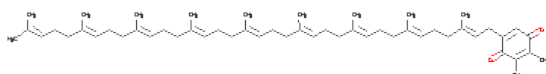
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
9	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
9	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
9	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
9	N	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
9	N	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
9	N	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
9	P	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 10 is 8-HYDROXY-5,7-DIMETHOXY-3-METHYL-2-TRIDECYL-4H-CHROME N-4-ONE (three-letter code: TDS) (formula: C<sub>25</sub>H<sub>38</sub>O<sub>5</sub>).



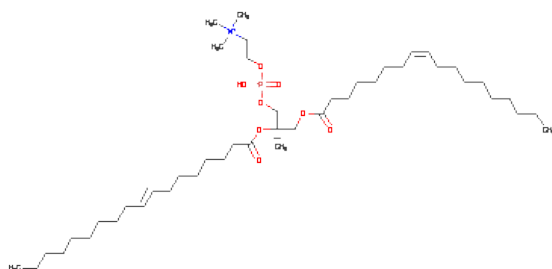
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			30	25	5		
10	N	1	Total	C	O	0	0
			30	25	5		

- Molecule 11 is 2,3-DIMETHYL-5-(3,7,11,15,19,23,27,31,35-NONAMETHYL-2,6,10,14,18,22,26,30,34-HEXATRIACONTANONAENYL-2,5-CYCLOHEXADIENE-1,4-DIONE-2,3-DIMETHYL-5-SOLANESYL-1,4-BENZOQUINONE (three-letter code: PL9) (formula:  $C_{53}H_{80}O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	A	1	Total	C	O	0	0
			55	53	2		
11	Q	1	Total	C	O	0	0
			55	53	2		

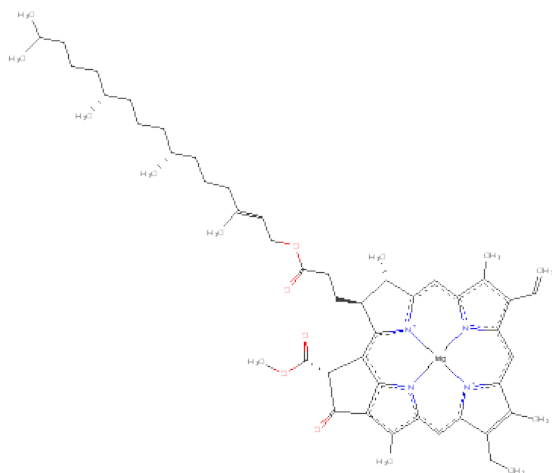
- Molecule 12 is (7R,17E)-4-HYDROXY-N,N,N,7-TETRAMETHYL-7-[(8E)-OCTADEC-8-ENOYLOXY]-10-OXO-3,5,9-TRIOXA-4-PHOSPHAHEPTACOS-17-EN-1-AMINIUM4-OXIDE (three-letter code: OPC) (formula: C<sub>45</sub>H<sub>87</sub>NO<sub>8</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
12	D	1	Total	C	N	O	P	0	0
			54	44	1	8	1		
12	B	1	Total	C	N	O	P	0	0
			54	44	1	8	1		
12	N	1	Total	C	N	O	P	0	0
			54	44	1	8	1		
12	Q	1	Total	C	N	O	P	0	0
			54	44	1	8	1		

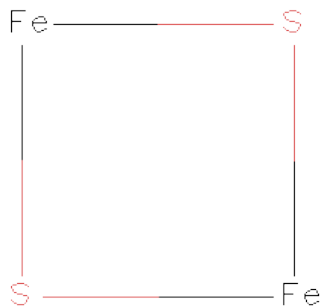
- Molecule 13 is CHLOROPHYLL A (three-letter code: CLA) (formula: C<sub>55</sub>H<sub>72</sub>MgN<sub>4</sub>O<sub>5</sub>).





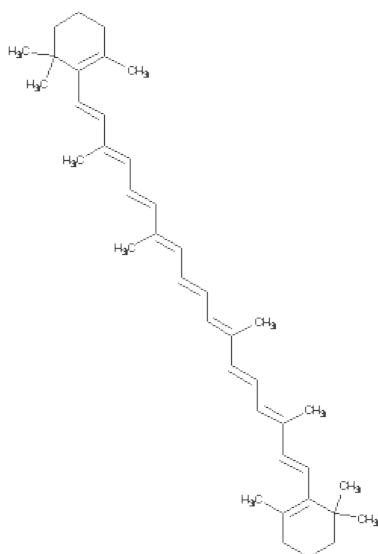
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
13	B	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
13	O	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		

- Molecule 14 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe<sub>2</sub>S<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
14	D	1	Total	Fe	S	0	0
			4	2	2		
14	Q	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 15 is BETA-CAROTENE (three-letter code: BCR) (formula:  $C_{40}H_{56}$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
15	E	1	Total C 40 40	0	0
15	R	1	Total C 40 40	0	0

- Molecule 16 is water.

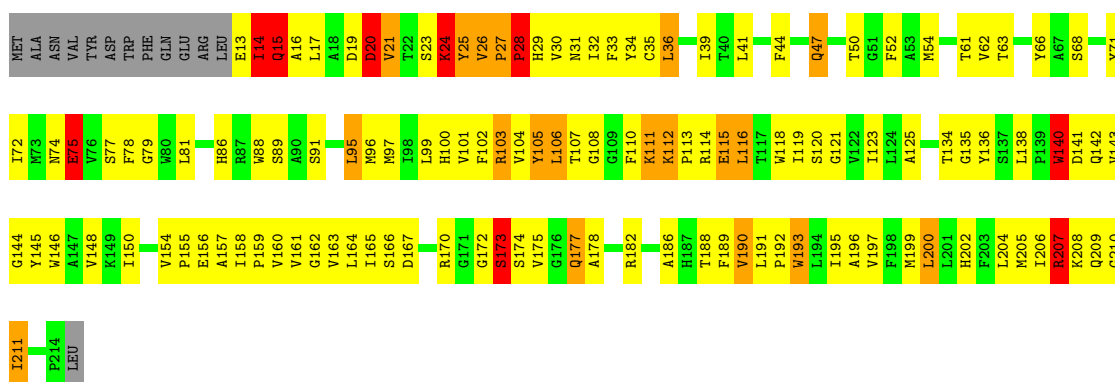
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
16	A	1	Total O 1 1	0	0
16	N	1	Total O 1 1	0	0

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

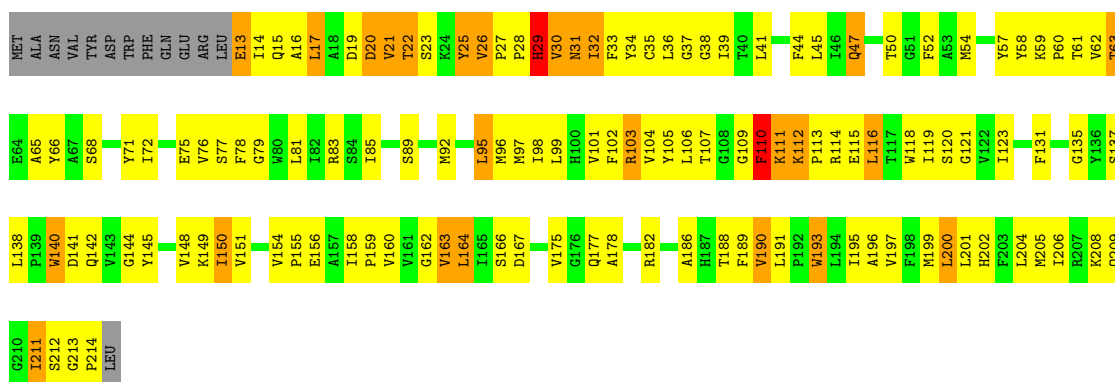
#### • Molecule 1: CYTOCHROME B6

Chain A:



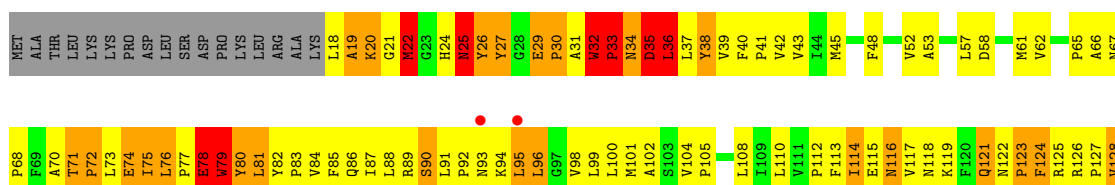
#### • Molecule 1: CYTOCHROME B6

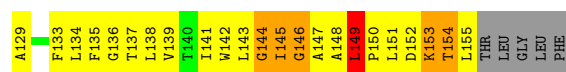
Chain N:



#### • Molecule 2: SUBUNIT IV

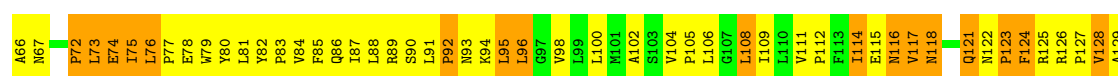
Chain B:





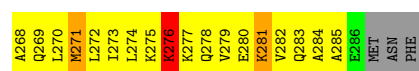
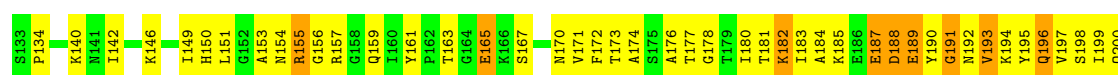
- Molecule 2: SUBUNIT IV

Chain O:



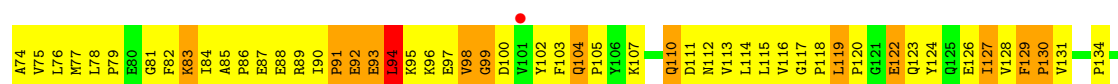
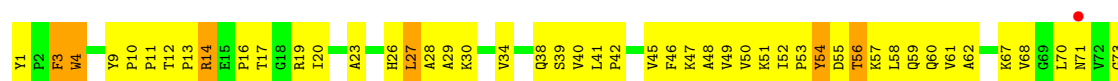
- Molecule 3: CYTOCHROME F

Chain C:



- Molecule 3: CYTOCHROME F

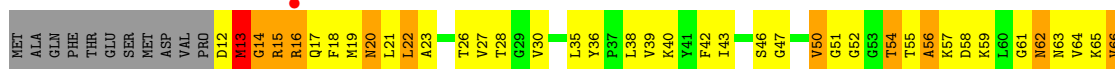
Chain P:





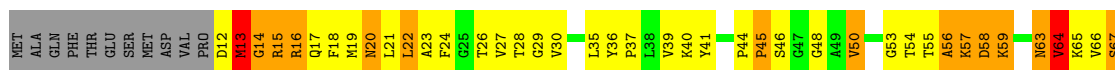
• Molecule 4: RIESKE IRON-SULFUR PROTEIN

Chain D:



• Molecule 4: RIESKE IRON-SULFUR PROTEIN

Chain Q:



• Molecule 5: PROTEIN PET L

Chain E:



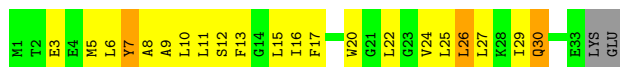
• Molecule 5: PROTEIN PET L

Chain R:



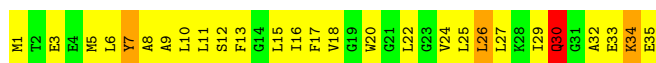
• Molecule 6: PROTEIN PET M

Chain F:



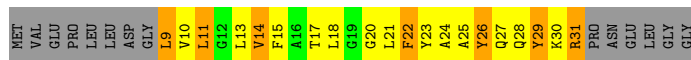
• Molecule 6: PROTEIN PET M

Chain S:



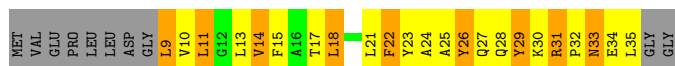
- Molecule 7: PROTEIN PET G

Chain G:



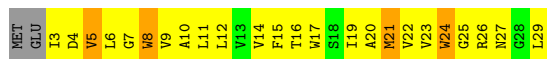
- Molecule 7: PROTEIN PET G

Chain T:



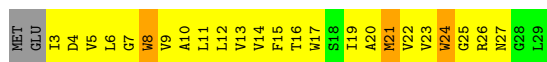
- Molecule 8: PROTEIN PET N

Chain H:



- Molecule 8: PROTEIN PET N

Chain U:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	157.54Å 157.54Å 360.35Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.16 – 3.00 48.16 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.16-3.00) 99.8 (48.16-3.00)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.30 (at 3.01Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.258 , 0.346 0.259 , 0.349	Depositor DCC
$R_{free}$ test set	2788 reflections (2.85%)	DCC
Wilson B-factor (Å <sup>2</sup> )	76.0	Xtriage
Anisotropy	0.238	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 110.1	EDS
Estimated twinning fraction	0.499 for h,-h-k,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 100543 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	15091	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.26% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities 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: CLA, PL9, FES, OPC, TDS, HEM, BCR

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	0.55	0/1641	0.77	2/2239 (0.1%)
1	N	0.56	0/1641	0.81	2/2239 (0.1%)
2	B	0.54	0/1110	0.88	3/1526 (0.2%)
2	O	0.58	0/1110	0.99	8/1526 (0.5%)
3	C	0.44	0/2248	0.73	1/3061 (0.0%)
3	P	0.45	0/2248	0.72	0/3061
4	D	0.47	0/1312	0.86	2/1786 (0.1%)
4	Q	0.47	0/1312	0.76	0/1786
5	E	0.67	0/253	0.80	0/340
5	R	0.64	0/253	0.77	0/340
6	F	0.58	0/255	0.66	0/343
6	S	0.52	0/274	0.61	0/366
7	G	0.64	0/188	0.91	0/253
7	T	0.63	0/221	0.95	0/299
8	H	0.61	0/220	0.83	1/301 (0.3%)
8	U	0.59	0/220	0.81	0/301
All	All	0.52	0/14506	0.80	19/19767 (0.1%)

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

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

There are no bond length outliers.

All (19) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	63	GLY	N-CA-C	7.67	132.29	113.10
2	B	36	LEU	N-CA-C	-7.09	91.85	111.00
2	O	64	GLU	N-CA-C	6.99	129.88	111.00
2	O	36	LEU	N-CA-C	-6.86	92.47	111.00
1	N	20	ASP	N-CA-C	6.33	128.08	111.00
2	O	117	VAL	N-CA-C	-6.17	94.34	111.00
1	A	207	ARG	N-CA-C	-5.99	94.84	111.00
2	B	144	GLY	N-CA-C	-5.87	98.42	113.10
2	O	144	GLY	N-CA-C	-5.82	98.55	113.10
2	B	35	ASP	N-CA-C	5.52	125.89	111.00
1	A	31	ASN	N-CA-C	5.51	125.89	111.00
2	O	35	ASP	N-CA-C	5.50	125.84	111.00
1	N	29	HIS	N-CA-C	5.40	125.59	111.00
4	D	54	THR	N-CA-C	5.38	125.52	111.00
3	C	231	LEU	CA-CB-CG	5.27	127.43	115.30
4	D	46	SER	N-CA-C	5.27	125.23	111.00
2	O	20	LYS	N-CA-C	5.16	124.94	111.00
2	O	39	VAL	N-CA-C	-5.01	97.47	111.00
8	H	5	VAL	N-CA-C	-5.00	97.49	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	N	105	TYR	Sidechain

## 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	1593	0	1623	242	0
1	N	1593	0	1623	230	0
2	B	1075	0	1117	258	0
2	O	1075	0	1117	284	0
3	C	2200	0	2216	355	0
3	P	2200	0	2216	338	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	1280	0	1263	317	0
4	Q	1280	0	1263	380	0
5	E	248	0	284	72	0
5	R	248	0	284	113	0
6	F	251	0	266	39	0
6	S	270	0	285	42	0
7	G	184	0	190	104	0
7	T	216	0	220	70	0
8	H	214	0	221	42	0
8	U	214	0	221	43	0
9	A	129	0	90	18	0
9	C	43	0	30	0	0
9	N	129	0	90	17	0
9	P	43	0	30	1	0
10	A	30	0	37	13	0
10	N	30	0	37	12	0
11	A	55	0	80	18	0
11	Q	55	0	80	23	0
12	B	54	0	83	14	0
12	D	54	0	83	24	0
12	N	54	0	83	10	0
12	Q	54	0	83	30	0
13	B	65	0	70	8	0
13	O	65	0	70	20	0
14	D	4	0	0	1	0
14	Q	4	0	0	1	0
15	E	40	0	56	7	0
15	R	40	0	56	7	0
16	A	1	0	0	0	0
16	N	1	0	0	0	0
All	All	15091	0	15467	2688	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 88.

All (2688) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:O:38:TYR:HB3	3:P:276:LYS:HG2	1.22	1.20
1:A:14:ILE:HA	1:A:17:LEU:HB2	1.26	1.18
1:N:214:PRO:HB3	5:R:29:ILE:HG22	1.19	1.17
4:D:166:THR:HA	4:D:179:VAL:HG13	1.22	1.17
2:B:71:THR:HB	2:B:72:PRO:HD3	1.28	1.15

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
11:A:305:PL9:H412	11:A:305:PL9:H462	1.18	1.13
2:B:20:LYS:HG3	2:B:21:GLY:H	1.05	1.13
11:Q:1305:PL9:H212	11:Q:1305:PL9:H262	1.24	1.12
12:D:306:OPC:HBI1	12:D:306:OPC:HAP2	1.31	1.12
4:Q:169:ASP:HB2	4:Q:176:PRO:HA	1.32	1.11
5:E:7:PHE:HA	5:E:10:VAL:HG12	1.29	1.11
12:Q:1307:OPC:HBE1	12:Q:1307:OPC:HBU2	1.11	1.11
4:Q:94:GLU:HB3	4:Q:100:ARG:HD3	1.18	1.10
4:D:136:THR:HB	4:D:138:ARG:HD2	1.25	1.10
4:Q:168:THR:HA	4:Q:176:PRO:HD3	1.19	1.10
4:D:124:PHE:HB2	4:D:133:TYR:HB2	1.32	1.09
2:B:126:ARG:HE	2:B:128:VAL:HG23	0.99	1.08
3:C:255:VAL:HG11	7:G:14:VAL:HG12	1.36	1.08
2:O:118:ASN:HD22	2:O:121:GLN:NE2	1.51	1.08
5:E:5:ALA:HB1	6:F:9:ALA:HB2	1.35	1.08
3:C:273:ILE:HD13	7:G:30:LYS:HD3	1.20	1.07
2:O:132:ILE:HG22	13:O:1201:CLA:HBB2	1.35	1.07
4:Q:156:GLN:HB2	4:Q:161:VAL:HG21	1.37	1.07
3:C:172:PHE:H	3:C:231:LEU:HG	1.17	1.07
5:R:7:PHE:HA	5:R:10:VAL:HG12	1.34	1.06
1:N:27:PRO:HG3	2:O:33:PRO:HG3	1.34	1.06
4:D:66:VAL:HG22	4:D:160:ILE:HG13	1.07	1.06
4:Q:76:GLY:H	4:Q:93:VAL:HG13	1.12	1.06
2:O:64:GLU:HG3	2:O:65:PRO:HD3	1.37	1.05
7:T:25:ALA:HB1	7:T:29:TYR:CZ	1.89	1.05
1:A:209:GLN:HG2	1:A:210:GLY:H	1.19	1.05
5:R:5:ALA:HB1	6:S:9:ALA:HB2	1.30	1.05
11:A:305:PL9:H203	12:D:306:OPC:HBE2	1.37	1.05
5:E:16:PHE:HB3	15:E:101:BCR:H363	1.36	1.05
12:N:1306:OPC:HAP2	12:N:1306:OPC:HBI1	1.39	1.05
3:C:281:LYS:HA	3:C:284:ALA:HB3	1.34	1.04
4:D:70:LEU:HD12	4:D:160:ILE:HD11	1.34	1.04
1:N:29:HIS:O	1:N:30:VAL:HG13	1.56	1.04
4:Q:56:ALA:HA	4:Q:81:VAL:HG13	1.36	1.03
3:P:281:LYS:HA	3:P:284:ALA:HB3	1.37	1.03
3:C:278:GLN:HE21	7:G:30:LYS:HE2	1.20	1.03
4:D:155:VAL:HG22	4:D:160:ILE:HG12	1.39	1.03
3:C:93:GLU:HA	3:C:97:GLU:HB2	1.38	1.02
1:A:207:ARG:HB2	1:A:207:ARG:HH11	1.21	1.02
4:D:51:GLY:HA2	4:D:164:PRO:HG2	1.35	1.02
4:Q:155:VAL:HG12	4:Q:157:ASP:H	1.21	1.01
2:B:18:LEU:HD13	2:B:31:ALA:HB2	1.42	1.01

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
7:G:25:ALA:HB1	7:G:29:TYR:CZ	1.95	1.01
3:C:172:PHE:O	3:C:231:LEU:HB3	1.60	1.01
5:R:16:PHE:HB3	15:R:1101:BCR:H363	1.38	1.01
2:O:124:PHE:CZ	5:R:26:ILE:HB	1.95	1.01
4:Q:116:PRO:HD2	4:Q:126:CYS:HA	1.41	1.00
3:C:271:MET:HG3	4:D:22:LEU:HD21	1.37	1.00
3:C:28:ALA:HB2	3:C:236:ASN:HD21	1.24	1.00
3:C:262:ILE:HG13	7:G:21:LEU:HG	1.38	1.00
7:G:26:TYR:HA	7:G:29:TYR:CD1	1.97	1.00
5:R:9:ILE:HG21	6:S:9:ALA:O	1.62	1.00
4:D:123:LYS:HB3	4:D:133:TYR:O	1.62	0.99
3:P:94:LEU:HD22	3:P:94:LEU:H	1.28	0.99
2:B:20:LYS:HG3	2:B:21:GLY:N	1.70	0.99
4:D:116:PRO:HG3	4:D:127:PRO:HD3	1.40	0.99
2:B:31:ALA:O	2:B:32:TRP:HB2	1.59	0.98
1:N:61:THR:HG22	1:N:63:THR:H	1.26	0.98
4:Q:57:LYS:HG3	4:Q:82:GLN:HE21	1.27	0.98
7:T:23:TYR:HA	7:T:26:TYR:CG	1.98	0.98
3:P:28:ALA:HB2	3:P:236:ASN:HD21	1.29	0.98
1:A:29:HIS:NE2	1:A:30:VAL:HG22	1.79	0.97
2:B:149:LEU:CD1	2:B:150:PRO:HD2	1.94	0.97
3:P:202:ASP:HA	3:P:206:THR:HB	1.43	0.97
4:D:51:GLY:HA2	4:D:164:PRO:CG	1.95	0.97
3:P:146:LYS:HD3	3:P:246:GLU:HG3	1.46	0.97
3:C:218:ILE:H	3:C:232:THR:HB	1.28	0.97
3:C:202:ASP:HA	3:C:206:THR:HB	1.45	0.97
2:B:61:MET:HB2	3:C:146:LYS:HD2	1.47	0.97
1:A:207:ARG:HB2	1:A:207:ARG:NH1	1.79	0.96
3:C:185:LYS:HG2	3:C:195:TYR:HB3	1.43	0.96
3:P:93:GLU:HA	3:P:97:GLU:HB2	1.43	0.96
4:D:57:LYS:HD2	4:D:82:GLN:HE21	1.27	0.96
5:E:9:ILE:HG21	6:F:9:ALA:O	1.64	0.95
2:B:126:ARG:NE	2:B:128:VAL:HG23	1.82	0.95
3:P:185:LYS:HG2	3:P:195:TYR:HB3	1.47	0.95
1:N:112:LYS:HB2	1:N:113:PRO:HD3	1.49	0.95
4:Q:123:LYS:HD3	4:Q:140:ILE:HD12	1.45	0.94
7:G:23:TYR:HA	7:G:26:TYR:CG	2.02	0.94
4:Q:74:ASN:H	4:Q:93:VAL:HG11	1.30	0.94
12:D:306:OPC:HBL1	12:D:306:OPC:HAR2	1.46	0.94
1:N:154:VAL:HG21	10:N:1304:TDS:HAY2	1.50	0.94
1:N:66:TYR:HB2	2:O:65:PRO:HG2	1.49	0.94
2:O:126:ARG:HH11	2:O:126:ARG:HG2	1.33	0.94

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:D:166:THR:HA	4:D:179:VAL:CG1	1.98	0.93
1:N:95:LEU:HD11	5:R:10:VAL:HG11	1.50	0.93
2:O:118:ASN:HD22	2:O:121:GLN:HE22	1.17	0.93
1:A:161:VAL:O	1:A:165:ILE:HG12	1.68	0.92
3:P:168:ASN:HA	3:P:172:PHE:CZ	2.04	0.92
2:O:126:ARG:NH1	2:O:128:VAL:HG23	1.83	0.92
3:P:172:PHE:H	3:P:231:LEU:HG	1.33	0.92
4:Q:73:HIS:CD2	4:Q:77:ASP:HB2	2.05	0.92
3:C:94:LEU:HD22	3:C:94:LEU:H	1.34	0.91
3:C:146:LYS:HD3	3:C:246:GLU:HG3	1.50	0.91
1:N:142:GLN:HG2	2:O:64:GLU:HG2	1.51	0.91
3:P:218:ILE:H	3:P:232:THR:HB	1.34	0.91
3:C:60:GLN:HE22	3:C:157:ARG:HG3	1.35	0.91
8:U:4:ASP:O	8:U:8:TRP:HB2	1.70	0.91
1:A:209:GLN:CG	1:A:210:GLY:H	1.79	0.90
3:P:172:PHE:O	3:P:231:LEU:HB3	1.71	0.90
1:A:170:ARG:HD3	1:A:174:SER:O	1.68	0.90
2:O:124:PHE:HZ	5:R:26:ILE:HB	1.36	0.90
2:O:31:ALA:O	2:O:32:TRP:HB2	1.70	0.90
1:A:114:ARG:HH12	1:A:209:GLN:H	1.06	0.90
4:D:99:ILE:HB	4:D:153:ALA:HB1	1.53	0.90
1:A:211:ILE:HD12	1:A:211:ILE:H	1.37	0.90
4:D:81:VAL:HG12	4:D:82:GLN:H	1.36	0.90
7:T:26:TYR:HA	7:T:29:TYR:CD1	2.06	0.90
1:A:209:GLN:HG2	1:A:210:GLY:N	1.87	0.90
4:Q:76:GLY:N	4:Q:93:VAL:HG13	1.86	0.89
1:A:24:LYS:HD3	1:A:24:LYS:H	1.35	0.89
2:O:80:TYR:HB3	13:O:1201:CLA:HED1	1.52	0.89
1:A:190:VAL:HG11	10:A:304:TDS:HAA3	1.54	0.89
4:Q:82:GLN:N	4:Q:89:THR:OG1	2.05	0.89
2:O:105:PRO:HG3	13:O:1201:CLA:H112	1.54	0.89
4:D:66:VAL:CG2	4:D:160:ILE:HG13	2.00	0.89
4:Q:57:LYS:HG3	4:Q:82:GLN:NE2	1.88	0.89
11:A:305:PL9:C46	11:A:305:PL9:H412	2.01	0.89
4:Q:91:ILE:HD12	4:Q:160:ILE:HD12	1.54	0.88
1:N:213:GLY:O	5:R:30:LYS:HD2	1.73	0.88
3:P:275:LYS:CE	4:Q:20:ASN:HB3	2.04	0.88
4:Q:146:LEU:HD13	4:Q:177:TRP:HB2	1.56	0.88
3:P:83:LYS:HB2	3:P:111:ASP:HB3	1.56	0.88
3:C:280:GLU:HG2	3:C:281:LYS:HZ3	1.39	0.87
3:P:231:LEU:HD13	3:P:232:THR:H	1.37	0.87
3:C:278:GLN:NE2	7:G:30:LYS:HE2	1.90	0.87

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:C:226:LYS:NZ	3:C:230:ALA:HB3	1.88	0.87
4:Q:78:ARG:HH22	4:Q:100:ARG:NH2	1.71	0.87
3:P:60:GLN:HE22	3:P:157:ARG:HG3	1.40	0.87
4:Q:156:GLN:HB2	4:Q:161:VAL:CG2	2.04	0.87
2:O:79:TRP:CZ2	5:R:1:MET:HB2	2.09	0.87
4:D:102:TYR:HB2	4:D:150:LEU:HG	1.56	0.87
7:T:21:LEU:HD22	7:T:21:LEU:O	1.75	0.87
1:N:110:PHE:H	1:N:110:PHE:HD2	1.20	0.87
8:H:4:ASP:O	8:H:8:TRP:HB2	1.74	0.87
3:P:226:LYS:NZ	3:P:230:ALA:HB3	1.90	0.87
3:C:255:VAL:CG1	7:G:14:VAL:HG12	2.04	0.86
3:P:275:LYS:HE3	4:Q:20:ASN:HB3	1.55	0.86
4:Q:77:ASP:H	4:Q:93:VAL:HG12	1.40	0.86
4:D:117:TRP:HA	4:D:124:PHE:HD1	1.41	0.86
2:O:126:ARG:HG2	2:O:126:ARG:NH1	1.89	0.86
1:N:103:ARG:HE	1:N:107:THR:HG21	1.39	0.86
4:Q:163:THR:OG1	4:Q:164:PRO:HD2	1.73	0.86
4:Q:73:HIS:CG	4:Q:77:ASP:HB2	2.10	0.86
2:O:64:GLU:HG3	2:O:65:PRO:CD	2.05	0.86
2:B:38:TYR:HB3	3:C:276:LYS:HG2	1.58	0.86
1:A:159:PRO:O	1:A:161:VAL:N	2.08	0.86
2:B:149:LEU:HD22	4:Q:129:HIS:HA	1.57	0.86
4:Q:22:LEU:HD23	4:Q:23:ALA:N	1.91	0.86
1:A:101:VAL:HG11	13:B:201:CLA:HMA1	1.58	0.85
12:N:1306:OPC:HBY1	12:N:1306:OPC:HBF2	1.56	0.85
4:Q:133:TYR:HA	4:Q:139:VAL:HA	1.57	0.85
1:A:211:ILE:HG22	5:E:26:ILE:HG12	1.59	0.85
4:D:142:GLY:HA2	4:D:144:ALA:H	1.42	0.85
4:D:22:LEU:HD23	4:D:23:ALA:N	1.92	0.85
1:A:158:ILE:HG23	1:A:159:PRO:HD2	1.59	0.85
1:A:170:ARG:HD2	1:A:172:GLY:O	1.76	0.85
3:C:280:GLU:HG2	3:C:281:LYS:NZ	1.91	0.85
2:B:149:LEU:CD2	4:Q:129:HIS:HA	2.06	0.85
1:A:14:ILE:HD13	1:A:15:GLN:N	1.92	0.85
4:Q:138:ARG:HE	4:Q:171:ARG:HD3	1.43	0.84
5:E:7:PHE:HA	5:E:10:VAL:CG1	2.07	0.84
2:O:149:LEU:O	2:O:151:LEU:N	2.09	0.84
1:A:14:ILE:HA	1:A:17:LEU:CB	2.07	0.84
2:O:38:TYR:CB	3:P:276:LYS:HG2	2.07	0.84
4:Q:105:ASN:HB3	4:Q:149:ALA:HB3	1.58	0.84
1:N:155:PRO:HG2	1:N:166:SER:OG	1.78	0.84
2:B:149:LEU:HD12	2:B:150:PRO:HD2	1.58	0.83

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:142:GLN:HE21	2:B:66:ALA:HA	1.42	0.83
4:Q:73:HIS:HB3	4:Q:93:VAL:HG11	1.57	0.83
3:C:83:LYS:HB2	3:C:111:ASP:HB3	1.58	0.83
4:D:142:GLY:HA2	4:D:144:ALA:N	1.94	0.83
4:Q:150:LEU:H	4:Q:178:TRP:HZ2	1.25	0.83
4:Q:64:VAL:HG12	4:Q:91:ILE:CG1	2.08	0.83
3:P:184:ALA:HB2	3:P:198:SER:OG	1.79	0.83
2:O:122:ASN:HD21	5:R:30:LYS:HD3	1.44	0.83
2:O:79:TRP:HZ2	5:R:1:MET:H3	1.27	0.83
3:C:218:ILE:N	3:C:232:THR:HB	1.94	0.83
3:C:26:HIS:CG	3:C:154:ASN:HD21	1.97	0.83
4:D:165:TRP:HD1	4:D:179:VAL:HA	1.44	0.82
4:Q:176:PRO:HG2	4:Q:179:VAL:HB	1.58	0.82
3:C:255:VAL:HG11	7:G:14:VAL:CG1	2.08	0.82
2:O:67:ASN:ND2	3:P:16:PRO:HB3	1.94	0.82
4:Q:78:ARG:HH22	4:Q:100:ARG:HH22	1.26	0.82
11:Q:1305:PL9:H302	12:Q:1307:OPC:HBC1	1.60	0.82
1:A:14:ILE:HD13	1:A:15:GLN:H	1.45	0.82
3:P:275:LYS:NZ	4:Q:20:ASN:HD22	1.77	0.82
12:B:307:OPC:HBI2	12:B:307:OPC:HAP1	1.59	0.82
4:D:154:THR:O	4:D:160:ILE:HG23	1.78	0.82
4:Q:94:GLU:CB	4:Q:100:ARG:HD3	2.05	0.82
3:P:168:ASN:HA	3:P:172:PHE:HZ	1.39	0.82
7:G:27:GLN:O	7:G:31:ARG:NH1	2.13	0.82
2:B:137:THR:O	2:B:141:ILE:HG12	1.80	0.82
3:C:226:LYS:HZ2	3:C:230:ALA:HB3	1.43	0.81
12:D:306:OPC:HBI1	12:D:306:OPC:CAP	2.08	0.81
2:O:126:ARG:CZ	2:O:128:VAL:HG23	2.10	0.81
1:A:62:VAL:HG12	1:A:140:TRP:CZ2	2.16	0.81
1:A:44:PHE:HD1	9:A:301:HEM:HBB1	1.45	0.81
4:Q:100:ARG:HD2	4:Q:102:TYR:OH	1.80	0.81
2:O:95:LEU:N	2:O:95:LEU:HD23	1.96	0.81
3:C:171:VAL:CG1	3:C:234:ASN:HB2	2.10	0.81
12:Q:1307:OPC:HBF2	12:Q:1307:OPC:HBX2	1.63	0.81
3:P:231:LEU:O	3:P:232:THR:HG23	1.80	0.81
3:P:26:HIS:CG	3:P:154:ASN:HD21	1.98	0.81
3:C:183:ILE:HD12	3:C:183:ILE:O	1.81	0.81
4:Q:155:VAL:HA	4:Q:161:VAL:H	1.43	0.81
3:C:93:GLU:HA	3:C:97:GLU:CB	2.11	0.81
3:P:90:ILE:HG23	3:P:95:LYS:HB2	1.62	0.81
3:P:217:LEU:HA	3:P:232:THR:HB	1.63	0.81
1:A:112:LYS:HA	1:A:115:GLU:OE1	1.80	0.81

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:14:ILE:CA	1:A:17:LEU:HB2	2.11	0.80
1:N:214:PRO:HB3	5:R:29:ILE:CG2	2.09	0.80
2:O:121:GLN:HB2	2:O:125:ARG:HD2	1.63	0.80
12:Q:1307:OPC:HAP1	12:Q:1307:OPC:OBJ	1.80	0.80
11:Q:1305:PL9:H462	11:Q:1305:PL9:H412	1.61	0.80
3:P:161:TYR:HE1	3:P:167:SER:HA	1.47	0.80
4:D:116:PRO:HG3	4:D:127:PRO:CD	2.11	0.80
11:Q:1305:PL9:H212	11:Q:1305:PL9:C26	2.10	0.80
7:T:11:LEU:H	7:T:11:LEU:HD22	1.44	0.80
3:P:183:ILE:O	3:P:183:ILE:HD12	1.80	0.80
3:C:269:GLN:HE22	7:G:24:ALA:CB	1.94	0.80
3:C:184:ALA:HB2	3:C:198:SER:OG	1.81	0.80
3:P:280:GLU:HG2	3:P:281:LYS:NZ	1.96	0.80
3:C:90:ILE:HG23	3:C:95:LYS:HB2	1.64	0.80
12:N:1306:OPC:HBO1	12:N:1306:OPC:HAT1	1.64	0.80
1:A:155:PRO:HG2	1:A:166:SER:OG	1.81	0.80
7:G:30:LYS:HB3	7:G:31:ARG:NH2	1.95	0.80
4:D:138:ARG:HB3	4:D:138:ARG:NH1	1.97	0.80
7:G:30:LYS:O	7:G:31:ARG:NH2	2.15	0.80
4:D:58:ASP:HB2	4:D:62:ASN:O	1.82	0.79
4:Q:139:VAL:HG21	4:Q:144:ALA:HB3	1.64	0.79
3:C:231:LEU:O	3:C:232:THR:HG23	1.82	0.79
3:C:199:ILE:O	3:C:208:VAL:HG12	1.81	0.79
1:A:206:ILE:C	1:A:207:ARG:HD3	2.03	0.79
3:C:91:PRO:HB2	3:C:94:LEU:HB2	1.65	0.79
7:T:23:TYR:HA	7:T:26:TYR:CD1	2.18	0.79
3:C:161:TYR:HE1	3:C:167:SER:HA	1.48	0.79
2:O:122:ASN:HD21	5:R:30:LYS:CD	1.94	0.79
1:A:114:ARG:HH12	1:A:209:GLN:N	1.81	0.79
7:G:11:LEU:H	7:G:11:LEU:HD22	1.44	0.79
2:O:122:ASN:ND2	5:R:30:LYS:HD3	1.98	0.79
4:D:93:VAL:HA	4:D:100:ARG:HB2	1.64	0.79
4:D:156:GLN:OE1	4:D:161:VAL:HG21	1.83	0.79
11:A:305:PL9:H101	12:D:306:OPC:HBF2	1.65	0.79
3:C:102:TYR:H	3:C:118:PRO:HG3	1.46	0.79
12:D:306:OPC:HAP2	12:D:306:OPC:CBI	2.07	0.78
4:D:70:LEU:HD12	4:D:160:ILE:CD1	2.13	0.78
2:O:57:LEU:HD22	7:T:14:VAL:CG2	2.13	0.78
2:B:18:LEU:HD13	2:B:31:ALA:CB	2.14	0.78
5:E:20:VAL:HG12	5:E:21:GLY:N	1.99	0.78
11:Q:1305:PL9:H321	12:Q:1307:OPC:HBW2	1.66	0.78
3:P:219:VAL:HG11	3:P:223:GLN:HE21	1.48	0.78

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:D:64:VAL:HG11	4:D:81:VAL:HG13	1.66	0.78
1:N:71:TYR:HA	1:N:75:GLU:OE1	1.84	0.78
7:T:23:TYR:HA	7:T:26:TYR:CD2	2.17	0.78
2:B:71:THR:HB	2:B:72:PRO:CD	2.11	0.78
1:A:29:HIS:CD2	1:A:30:VAL:HG13	2.19	0.78
1:N:195:ILE:O	1:N:199:MET:HG3	1.84	0.78
2:O:112:PRO:O	2:O:116:ASN:HB2	1.84	0.78
1:N:114:ARG:HH21	1:N:209:GLN:N	1.82	0.77
2:O:149:LEU:HB3	2:O:150:PRO:HD2	1.66	0.77
3:P:134:PRO:HB3	3:P:142:ILE:HD13	1.65	0.77
4:Q:154:THR:O	4:Q:160:ILE:HA	1.83	0.77
1:A:28:PRO:HD2	2:B:31:ALA:O	1.85	0.77
3:P:188:ASP:CB	3:P:193:VAL:HA	2.15	0.77
1:A:112:LYS:H	1:A:113:PRO:HD2	1.50	0.77
4:Q:169:ASP:HB2	4:Q:176:PRO:CA	2.12	0.77
1:A:39:ILE:HD11	15:E:101:BCR:H312	1.65	0.77
12:Q:1307:OPC:HAL2	12:Q:1307:OPC:CAP	2.15	0.77
3:P:255:VAL:HG11	7:T:14:VAL:HG12	1.66	0.77
4:Q:155:VAL:HG12	4:Q:157:ASP:N	1.98	0.77
8:H:19:ILE:O	8:H:23:VAL:HG22	1.84	0.77
7:G:21:LEU:O	7:G:21:LEU:HD22	1.84	0.77
2:O:149:LEU:HB3	2:O:150:PRO:CD	2.14	0.77
3:C:231:LEU:HD13	3:C:232:THR:H	1.49	0.77
1:N:116:LEU:H	1:N:116:LEU:HD12	1.50	0.77
5:R:23:ILE:O	5:R:27:LYS:HB2	1.85	0.77
4:Q:176:PRO:CG	4:Q:179:VAL:HB	2.14	0.77
3:C:92:GLU:O	3:C:96:LYS:HB3	1.85	0.77
3:P:102:TYR:H	3:P:118:PRO:HG3	1.49	0.77
3:C:259:ILE:HA	7:G:17:THR:HG21	1.65	0.77
2:O:38:TYR:CD1	3:P:276:LYS:HD3	2.20	0.77
2:B:34:ASN:HD21	12:B:307:OPC:CBK	1.97	0.77
12:Q:1307:OPC:HBU2	12:Q:1307:OPC:CBE	2.05	0.76
1:A:47:GLN:HE22	1:A:89:SER:CB	1.97	0.76
8:U:19:ILE:O	8:U:23:VAL:HG22	1.85	0.76
3:C:188:ASP:CB	3:C:193:VAL:HA	2.14	0.76
3:C:28:ALA:HB2	3:C:236:ASN:ND2	1.98	0.76
2:B:95:LEU:HD23	2:B:95:LEU:N	2.00	0.76
1:N:29:HIS:N	1:N:29:HIS:ND1	2.31	0.76
4:D:57:LYS:HB2	4:D:82:GLN:HG2	1.66	0.76
3:C:281:LYS:CA	3:C:284:ALA:HB3	2.16	0.76
7:G:25:ALA:HB1	7:G:29:TYR:CE2	2.21	0.76
3:P:199:ILE:O	3:P:208:VAL:HG12	1.85	0.76

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:Q:150:LEU:HB2	4:Q:178:TRP:HH2	1.51	0.76
5:R:7:PHE:HA	5:R:10:VAL:CG1	2.12	0.76
2:O:53:ALA:O	2:O:57:LEU:HG	1.84	0.76
2:B:112:PRO:O	2:B:116:ASN:HB2	1.85	0.76
3:P:271:MET:HG3	4:Q:22:LEU:HD21	1.67	0.76
12:D:306:OPC:HBW2	12:D:306:OPC:HCA2	1.67	0.76
1:A:158:ILE:HB	1:A:162:GLY:HA3	1.66	0.76
2:O:79:TRP:CZ3	5:R:4:GLY:CA	2.69	0.76
4:Q:175:LYS:HB2	4:Q:179:VAL:HG11	1.66	0.76
3:C:269:GLN:HE22	7:G:24:ALA:HB1	1.50	0.76
2:B:151:LEU:HD12	2:B:153:LYS:HD3	1.67	0.76
5:E:23:ILE:O	5:E:27:LYS:HB2	1.86	0.76
4:Q:64:VAL:HG12	4:Q:91:ILE:HG12	1.64	0.76
3:P:218:ILE:N	3:P:232:THR:HB	1.99	0.76
3:C:181:THR:O	3:C:221:GLU:HB3	1.86	0.76
1:N:142:GLN:HG2	2:O:64:GLU:CG	2.15	0.75
1:N:141:ASP:HA	2:O:65:PRO:HD3	1.68	0.75
2:B:151:LEU:HA	2:B:153:LYS:HD3	1.67	0.75
3:P:231:LEU:HD11	3:P:233:ASN:O	1.85	0.75
4:D:117:TRP:HA	4:D:124:PHE:CD1	2.21	0.75
3:C:218:ILE:H	3:C:232:THR:CB	1.99	0.75
2:O:57:LEU:HD22	7:T:14:VAL:HG23	1.67	0.75
4:Q:64:VAL:HB	4:Q:81:VAL:HG22	1.67	0.75
1:A:195:ILE:O	1:A:199:MET:HG3	1.87	0.75
2:O:122:ASN:HD22	2:O:125:ARG:CZ	1.99	0.75
4:D:123:LYS:HA	4:D:134:ASP:C	2.07	0.75
2:O:79:TRP:HZ2	5:R:1:MET:HB2	1.48	0.75
3:P:91:PRO:HB2	3:P:94:LEU:HB2	1.66	0.75
1:A:158:ILE:HB	1:A:162:GLY:CA	2.16	0.75
4:D:84:LEU:HD22	4:D:87:ASP:HB2	1.67	0.75
4:Q:151:CYS:HB3	4:Q:162:LEU:HD23	1.69	0.75
2:B:122:ASN:ND2	2:B:124:PHE:HB3	2.00	0.75
7:T:25:ALA:CB	7:T:29:TYR:CZ	2.69	0.75
2:O:122:ASN:OD1	5:R:30:LYS:HD3	1.87	0.75
12:D:306:OPC:CBL	12:D:306:OPC:HAR2	2.17	0.75
7:G:23:TYR:HA	7:G:26:TYR:CD2	2.20	0.75
2:O:146:GLY:H	2:O:149:LEU:HD12	1.49	0.75
2:O:146:GLY:N	2:O:149:LEU:HD12	2.02	0.75
4:Q:156:GLN:O	4:Q:157:ASP:HB3	1.86	0.75
1:A:24:LYS:CD	1:A:24:LYS:H	2.00	0.75
3:C:39:SER:HB3	3:C:248:VAL:HB	1.66	0.75
4:D:81:VAL:HG21	4:D:91:ILE:HD11	1.68	0.75

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:Q:159:ASN:O	4:Q:161:VAL:HG13	1.85	0.75
4:Q:67:SER:C	4:Q:68:LYS:HD2	2.07	0.75
3:C:219:VAL:H	3:C:232:THR:HG22	1.52	0.75
5:E:16:PHE:HB3	15:E:101:BCR:C36	2.17	0.74
3:P:188:ASP:HB2	3:P:193:VAL:HA	1.69	0.74
2:O:117:VAL:HG13	2:O:118:ASN:OD1	1.87	0.74
2:B:122:ASN:HD22	2:B:124:PHE:HD2	1.35	0.74
1:A:47:GLN:HE22	1:A:89:SER:HB3	1.52	0.74
4:D:134:ASP:CG	4:D:135:GLU:H	1.90	0.74
4:D:65:LYS:HB3	4:D:68:LYS:HD3	1.68	0.74
4:Q:150:LEU:HD11	4:Q:171:ARG:HH21	1.53	0.74
4:Q:87:ASP:HB3	4:Q:88:PRO:HD2	1.67	0.74
4:Q:92:VAL:CG2	4:Q:100:ARG:HB2	2.17	0.74
3:P:52:ILE:HG12	3:P:153:ALA:HB1	1.69	0.74
4:Q:132:GLN:HE22	4:Q:141:ARG:HD3	1.53	0.74
11:A:305:PL9:H252	12:D:306:OPC:HBN1	1.68	0.74
3:P:229:GLU:HG3	3:P:230:ALA:H	1.53	0.74
10:A:304:TDS:HBC1	2:B:85:PHE:HA	1.70	0.73
3:C:182:LYS:HG2	3:C:198:SER:HB2	1.69	0.73
1:N:47:GLN:HE22	1:N:89:SER:HB3	1.53	0.73
2:O:76:LEU:HD12	5:R:1:MET:SD	2.28	0.73
1:A:29:HIS:CD2	1:A:30:VAL:HG22	2.21	0.73
7:G:25:ALA:CB	7:G:29:TYR:CZ	2.70	0.73
2:O:137:THR:O	2:O:141:ILE:HG12	1.89	0.73
2:O:79:TRP:CH2	5:R:4:GLY:HA3	2.24	0.73
3:P:28:ALA:HB2	3:P:236:ASN:ND2	2.03	0.73
1:A:146:TRP:CE3	2:B:71:THR:HG23	2.24	0.73
4:Q:155:VAL:C	4:Q:161:VAL:HG22	2.09	0.73
3:P:280:GLU:HG2	3:P:281:LYS:HZ3	1.52	0.73
2:B:79:TRP:HA	2:B:82:TYR:CE2	2.24	0.73
4:Q:20:ASN:C	4:Q:22:LEU:H	1.89	0.73
4:D:78:ARG:HE	4:D:92:VAL:CG1	2.02	0.73
2:B:133:PHE:O	2:B:137:THR:HG23	1.89	0.73
3:P:93:GLU:HA	3:P:97:GLU:CB	2.19	0.73
10:A:304:TDS:HBB2	2:B:88:LEU:HD22	1.70	0.73
2:B:144:GLY:O	2:B:145:ILE:HB	1.87	0.73
8:U:23:VAL:HA	8:U:26:ARG:HD2	1.71	0.73
4:D:133:TYR:HA	4:D:139:VAL:HA	1.71	0.73
4:D:138:ARG:HH21	4:D:171:ARG:CZ	2.02	0.73
4:Q:73:HIS:HB3	4:Q:93:VAL:CG1	2.18	0.73
11:A:305:PL9:H512	11:A:305:PL9:H401	1.71	0.73
3:C:188:ASP:HB2	3:C:193:VAL:HA	1.68	0.73

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:R:23:ILE:HG23	5:R:24:PHE:CD2	2.23	0.72
2:B:149:LEU:HD13	2:B:150:PRO:HD2	1.69	0.72
2:O:32:TRP:H	2:O:33:PRO:HD3	1.53	0.72
2:O:122:ASN:HD21	5:R:30:LYS:CG	2.01	0.72
4:D:165:TRP:CD1	4:D:179:VAL:HA	2.24	0.72
7:T:25:ALA:HB1	7:T:29:TYR:OH	1.89	0.72
2:B:118:ASN:HB3	2:B:121:GLN:O	1.89	0.72
3:C:229:GLU:HG3	3:C:230:ALA:H	1.54	0.72
4:Q:76:GLY:HA2	4:Q:93:VAL:O	1.88	0.72
2:B:32:TRP:HH2	3:C:277:LYS:HD2	1.55	0.72
12:Q:1307:OPC:HAQ1	12:Q:1307:OPC:HBG3	1.72	0.72
12:Q:1307:OPC:HBX2	12:Q:1307:OPC:CBF	2.19	0.72
3:C:217:LEU:HA	3:C:232:THR:HB	1.72	0.72
3:C:171:VAL:HG13	3:C:234:ASN:HB2	1.71	0.72
3:P:211:ILE:HD11	3:P:231:LEU:HB2	1.71	0.72
1:A:116:LEU:H	1:A:116:LEU:HD12	1.53	0.72
4:Q:131:SER:OG	4:Q:144:ALA:HB2	1.89	0.72
4:Q:91:ILE:HG23	4:Q:160:ILE:HD12	1.71	0.72
4:Q:178:TRP:HE3	4:Q:178:TRP:N	1.88	0.72
2:O:79:TRP:CZ3	5:R:4:GLY:HA3	2.25	0.72
4:D:66:VAL:HG21	4:D:155:VAL:HG13	1.71	0.72
6:S:6:LEU:O	6:S:10:LEU:HG	1.88	0.72
4:Q:103:GLY:O	4:Q:150:LEU:HA	1.89	0.72
5:E:23:ILE:HG23	5:E:24:PHE:CD2	2.25	0.72
1:N:200:LEU:O	1:N:204:LEU:HG	1.89	0.72
4:D:16:ARG:C	4:D:18:PHE:H	1.94	0.72
1:A:17:LEU:HD23	1:A:17:LEU:O	1.88	0.72
1:A:114:ARG:NH1	1:A:209:GLN:H	1.85	0.72
2:O:133:PHE:O	2:O:137:THR:HG23	1.90	0.72
1:A:135:GLY:O	9:A:301:HEM:HAA1	1.90	0.72
12:D:306:OPC:HBL1	12:D:306:OPC:HAP2	1.71	0.71
2:O:20:LYS:HD3	2:O:21:GLY:H	1.53	0.71
4:Q:137:GLY:O	4:Q:147:SER:HB3	1.90	0.71
4:Q:66:VAL:HG22	4:Q:160:ILE:CG1	2.20	0.71
3:C:278:GLN:NE2	6:F:30:GLN:OE1	2.23	0.71
3:P:219:VAL:H	3:P:232:THR:HG22	1.55	0.71
4:Q:105:ASN:HB3	4:Q:149:ALA:CB	2.20	0.71
8:H:11:LEU:HB3	8:H:15:PHE:CZ	2.24	0.71
1:A:103:ARG:HH11	1:A:103:ARG:HG3	1.54	0.71
3:P:218:ILE:H	3:P:232:THR:CB	2.03	0.71
6:F:6:LEU:O	6:F:10:LEU:HG	1.90	0.71
1:N:29:HIS:HE2	1:N:209:GLN:HG3	1.56	0.71

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:Q:64:VAL:HB	4:Q:81:VAL:CG2	2.20	0.71
1:N:66:TYR:CB	2:O:65:PRO:HG2	2.20	0.71
3:P:275:LYS:HZ3	3:P:276:LYS:NZ	1.89	0.71
3:P:275:LYS:HZ1	4:Q:20:ASN:HD22	1.36	0.71
2:O:106:LEU:HD21	13:O:1201:CLA:H151	1.72	0.71
3:P:185:LYS:HE2	3:P:195:TYR:HB3	1.71	0.71
2:O:142:TRP:O	2:O:145:ILE:HG12	1.91	0.71
2:O:84:VAL:O	2:O:88:LEU:HD13	1.90	0.71
7:T:14:VAL:O	7:T:17:THR:HB	1.91	0.71
4:D:57:LYS:CB	4:D:82:GLN:HG2	2.21	0.71
2:B:18:LEU:HD22	2:B:30:PRO:O	1.90	0.71
2:B:118:ASN:HB3	2:B:121:GLN:HE21	1.55	0.71
3:P:92:GLU:O	3:P:96:LYS:HB3	1.91	0.71
1:N:212:SER:OG	5:R:26:ILE:HG12	1.90	0.71
4:Q:117:TRP:HE1	4:Q:135:GLU:HA	1.55	0.71
2:B:126:ARG:HE	2:B:128:VAL:CG2	1.93	0.71
2:B:151:LEU:C	2:B:153:LYS:H	1.93	0.71
4:Q:145:PRO:HG2	4:Q:146:LEU:H	1.55	0.70
3:C:218:ILE:HG13	3:C:219:VAL:H	1.56	0.70
4:D:115:VAL:HG21	4:D:148:LEU:HD21	1.73	0.70
4:Q:56:ALA:O	4:Q:63:ASN:HA	1.91	0.70
1:A:211:ILE:HD11	2:B:122:ASN:OD1	1.91	0.70
4:D:15:ARG:NH2	4:D:17:GLN:HG2	2.05	0.70
6:F:5:MET:HA	6:F:8:ALA:HB3	1.73	0.70
7:T:11:LEU:O	7:T:14:VAL:HG22	1.90	0.70
2:O:132:ILE:CG2	13:O:1201:CLA:HBB2	2.18	0.70
3:C:226:LYS:HG2	3:C:229:GLU:HB3	1.73	0.70
3:P:3:PHE:HZ	3:P:119:LEU:HD11	1.56	0.70
4:D:147:SER:HB2	4:D:177:TRP:HH2	1.56	0.70
4:Q:64:VAL:HA	4:Q:69:PHE:CD1	2.26	0.70
11:Q:1305:PL9:C46	11:Q:1305:PL9:H412	2.21	0.70
3:C:219:VAL:HG11	3:C:223:GLN:HE21	1.54	0.70
3:P:172:PHE:CZ	3:P:212:PRO:HG3	2.25	0.70
1:A:111:LYS:HG2	1:A:112:LYS:N	2.06	0.70
1:N:189:PHE:O	1:N:193:TRP:HE3	1.74	0.70
1:N:209:GLN:NE2	2:O:28:GLY:H	1.90	0.70
1:A:29:HIS:HD2	1:A:30:VAL:HG13	1.56	0.70
4:D:20:ASN:C	4:D:22:LEU:H	1.91	0.70
1:A:189:PHE:O	1:A:193:TRP:HE3	1.75	0.70
4:Q:152:HIS:C	4:Q:162:LEU:HG	2.12	0.70
4:Q:93:VAL:CA	4:Q:100:ARG:HG3	2.21	0.70
2:B:145:ILE:O	2:B:146:GLY:C	2.29	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:O:117:VAL:O	2:O:118:ASN:HB2	1.92	0.70
6:S:5:MET:HA	6:S:8:ALA:HB3	1.73	0.70
2:B:57:LEU:HD21	3:C:258:MET:HE2	1.72	0.70
11:Q:1305:PL9:H462	11:Q:1305:PL9:C41	2.21	0.70
2:O:79:TRP:CZ3	5:R:4:GLY:HA2	2.26	0.70
2:B:144:GLY:O	2:B:145:ILE:CB	2.39	0.70
4:D:115:VAL:HG11	4:D:148:LEU:HD11	1.74	0.70
1:A:207:ARG:CB	1:A:207:ARG:HH11	2.00	0.70
2:O:79:TRP:CE3	5:R:4:GLY:HA2	2.26	0.70
3:C:232:THR:O	3:C:233:ASN:HB2	1.91	0.70
7:T:28:GLN:O	7:T:31:ARG:N	2.24	0.70
3:P:39:SER:HB3	3:P:248:VAL:HB	1.72	0.70
1:N:34:TYR:OH	1:N:211:ILE:HG22	1.92	0.69
4:D:168:THR:HA	4:D:176:PRO:HD3	1.74	0.69
2:B:20:LYS:NZ	2:B:21:GLY:O	2.23	0.69
1:N:21:VAL:HG12	1:N:22:THR:H	1.55	0.69
3:P:70:LEU:H	3:P:70:LEU:HD23	1.57	0.69
7:G:23:TYR:HA	7:G:26:TYR:CD1	2.26	0.69
5:E:7:PHE:CA	5:E:10:VAL:HG12	2.17	0.69
3:C:171:VAL:HB	3:C:231:LEU:HD12	1.75	0.69
2:B:26:TYR:N	2:B:26:TYR:HD2	1.90	0.69
2:O:122:ASN:ND2	5:R:26:ILE:HG22	2.08	0.69
4:D:57:LYS:HD2	4:D:82:GLN:NE2	2.04	0.69
1:A:26:VAL:HB	2:B:29:GLU:CB	2.22	0.69
3:P:182:LYS:HG2	3:P:198:SER:HB2	1.74	0.69
4:D:78:ARG:HE	4:D:92:VAL:HG11	1.58	0.69
4:D:57:LYS:CD	4:D:82:GLN:HE21	2.04	0.69
1:A:206:ILE:O	1:A:207:ARG:HD3	1.93	0.69
10:N:1304:TDS:HBB2	2:O:81:LEU:HD13	1.73	0.69
3:P:119:LEU:HD12	3:P:119:LEU:N	2.07	0.69
2:B:26:TYR:N	2:B:26:TYR:CD2	2.61	0.69
4:Q:138:ARG:HE	4:Q:171:ARG:CD	2.05	0.69
4:Q:69:PHE:O	4:Q:91:ILE:HG21	1.92	0.69
3:C:126:GLU:OE1	3:C:126:GLU:HA	1.92	0.69
4:Q:68:LYS:N	4:Q:68:LYS:HD2	2.08	0.69
1:N:111:LYS:HB2	1:N:114:ARG:HH11	1.58	0.69
2:O:122:ASN:HD22	2:O:125:ARG:NH2	1.90	0.69
4:D:132:GLN:HB2	4:D:140:ILE:HB	1.75	0.69
4:Q:165:TRP:HZ2	4:Q:176:PRO:HB3	1.54	0.69
2:B:127:PRO:C	2:B:129:ALA:H	1.94	0.69
2:B:22:MET:CE	2:B:22:MET:HA	2.23	0.69
3:C:172:PHE:CZ	3:C:212:PRO:HG3	2.28	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:C:172:PHE:C	3:C:231:LEU:HB3	2.13	0.69
3:C:119:LEU:HD12	3:C:119:LEU:N	2.08	0.69
6:S:25:LEU:HD23	6:S:25:LEU:O	1.92	0.69
5:R:18:ILE:O	5:R:22:ILE:HB	1.91	0.69
4:Q:123:LYS:HE3	4:Q:125:LYS:NZ	2.08	0.69
1:N:102:PHE:HB3	5:R:18:ILE:HD11	1.75	0.69
4:D:123:LYS:HA	4:D:134:ASP:O	1.91	0.69
4:Q:65:LYS:HE2	4:Q:158:ASP:O	1.90	0.69
7:G:11:LEU:HA	7:G:14:VAL:CG1	2.22	0.69
2:O:132:ILE:HG22	13:O:1201:CLA:CBB	2.19	0.69
3:P:178:GLY:HA2	3:P:225:VAL:HG13	1.75	0.69
1:N:110:PHE:N	1:N:110:PHE:HD2	1.90	0.68
5:R:20:VAL:HG12	5:R:21:GLY:N	2.06	0.68
2:B:18:LEU:HB3	2:B:30:PRO:O	1.93	0.68
7:G:14:VAL:CG2	7:G:15:PHE:N	2.56	0.68
12:Q:1307:OPC:HAX2	12:Q:1307:OPC:HAT1	1.76	0.68
2:O:74:GLU:HG2	2:O:75:ILE:N	2.08	0.68
3:C:277:LYS:HE3	7:G:31:ARG:HG3	1.74	0.68
4:D:103:GLY:O	4:D:151:CYS:HB2	1.92	0.68
4:D:154:THR:HG23	4:D:161:VAL:O	1.94	0.68
2:O:18:LEU:HD13	2:O:31:ALA:CB	2.23	0.68
4:Q:109:THR:O	4:Q:145:PRO:HG3	1.93	0.68
3:C:60:GLN:NE2	3:C:157:ARG:HG3	2.08	0.68
2:O:52:VAL:O	2:O:56:VAL:HG23	1.93	0.68
1:N:110:PHE:N	1:N:110:PHE:CD2	2.61	0.68
2:B:32:TRP:HZ3	2:B:36:LEU:HA	1.59	0.68
2:O:133:PHE:HA	13:O:1201:CLA:HMB1	1.75	0.68
3:P:161:TYR:HE1	3:P:167:SER:CA	2.07	0.68
3:C:134:PRO:HB3	3:C:142:ILE:HD13	1.75	0.68
4:Q:102:TYR:CE1	4:Q:136:THR:HG23	2.28	0.68
3:C:258:MET:O	3:C:262:ILE:HD13	1.93	0.68
4:Q:24:PHE:CE1	11:Q:1305:PL9:H202	2.28	0.68
3:C:180:ILE:HD11	3:C:182:LYS:O	1.93	0.68
4:D:126:CYS:HB3	14:D:200:FES:S1	2.34	0.68
7:T:25:ALA:HB1	7:T:29:TYR:CE2	2.27	0.68
2:O:127:PRO:C	2:O:129:ALA:H	1.95	0.68
3:P:226:LYS:HZ2	3:P:230:ALA:HB3	1.57	0.68
4:Q:92:VAL:HG23	4:Q:100:ARG:HB2	1.75	0.68
3:P:185:LYS:HD3	3:P:195:TYR:HD2	1.58	0.68
3:P:60:GLN:NE2	3:P:157:ARG:HG3	2.09	0.68
1:A:62:VAL:CG1	1:A:177:GLN:HB3	2.24	0.68
2:O:32:TRP:H	2:O:33:PRO:CD	2.06	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:Q:93:VAL:HA	4:Q:100:ARG:HG3	1.74	0.68
2:B:42:VAL:HG12	7:G:28:GLN:HE22	1.59	0.68
12:B:307:OPC:HBI2	12:B:307:OPC:CAP	2.24	0.68
3:P:181:THR:O	3:P:221:GLU:HB3	1.92	0.68
1:A:142:GLN:NE2	2:B:67:ASN:H	1.92	0.68
3:C:52:ILE:HG12	3:C:153:ALA:HB1	1.75	0.68
4:D:156:GLN:HB2	4:D:161:VAL:HG22	1.75	0.67
3:P:232:THR:O	3:P:233:ASN:HB2	1.92	0.67
7:G:11:LEU:O	7:G:14:VAL:HG22	1.94	0.67
1:N:101:VAL:HG11	13:O:1201:CLA:HMA2	1.76	0.67
4:D:117:TRP:CZ3	4:D:119:ALA:HA	2.30	0.67
4:Q:115:VAL:HG11	4:Q:124:PHE:HB3	1.76	0.67
6:F:26:LEU:CD2	8:H:17:TRP:HE1	2.07	0.67
2:O:89:ARG:HH12	2:O:148:ALA:H	1.43	0.67
7:T:14:VAL:CG2	7:T:15:PHE:N	2.56	0.67
8:U:11:LEU:HB3	8:U:15:PHE:CZ	2.29	0.67
1:N:188:THR:HG22	9:N:301:HEM:HBC2	1.75	0.67
1:N:149:LYS:HA	1:N:175:VAL:HG21	1.76	0.67
1:N:137:SER:O	1:N:140:TRP:HD1	1.78	0.67
3:C:1:TYR:HD2	3:C:119:LEU:HD21	1.59	0.67
3:C:269:GLN:OE1	7:G:24:ALA:HA	1.93	0.67
3:P:275:LYS:HE2	4:Q:17:GLN:HB2	1.76	0.67
4:Q:150:LEU:HB2	4:Q:178:TRP:CH2	2.29	0.67
4:Q:16:ARG:C	4:Q:18:PHE:H	1.98	0.67
11:A:305:PL9:H111	12:D:306:OPC:HBF2	1.75	0.67
2:O:79:TRP:HZ2	5:R:1:MET:CB	2.07	0.67
4:Q:81:VAL:HG12	4:Q:82:GLN:HG3	1.76	0.67
1:A:114:ARG:CZ	1:A:208:LYS:HG3	2.24	0.67
2:B:153:LYS:O	2:B:154:THR:O	2.12	0.67
1:N:110:PHE:HD1	2:O:112:PRO:HB3	1.60	0.67
4:Q:138:ARG:NE	4:Q:171:ARG:HD3	2.10	0.67
2:O:126:ARG:HH11	2:O:126:ARG:CG	2.04	0.67
4:D:94:GLU:OE2	4:D:98:ALA:HB3	1.95	0.67
4:Q:77:ASP:C	4:Q:78:ARG:HD3	2.15	0.67
5:R:26:ILE:HA	5:R:29:ILE:HG13	1.77	0.66
2:B:57:LEU:HD21	3:C:258:MET:CE	2.25	0.66
3:P:226:LYS:HZ1	3:P:230:ALA:HB3	1.60	0.66
2:B:38:TYR:OH	12:B:307:OPC:HAU2	1.94	0.66
7:G:9:LEU:N	7:G:9:LEU:HD23	2.10	0.66
3:C:211:ILE:HD11	3:C:231:LEU:HB2	1.76	0.66
3:P:171:VAL:HG13	3:P:234:ASN:HD22	1.61	0.66
3:C:226:LYS:HE3	3:C:229:GLU:HB3	1.77	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
8:U:4:ASP:O	8:U:8:TRP:CB	2.43	0.66
9:N:301:HEM:HBD1	9:N:301:HEM:HHA	1.75	0.66
4:Q:108:CYS:HB2	4:Q:113:CYS:O	1.95	0.66
4:Q:134:ASP:CG	4:Q:138:ARG:HB2	2.15	0.66
3:P:271:MET:HB3	4:Q:23:ALA:HA	1.77	0.66
4:Q:115:VAL:HG13	4:Q:126:CYS:HB2	1.78	0.66
2:O:149:LEU:C	2:O:151:LEU:H	1.98	0.66
6:S:34:LYS:HE2	6:S:35:GLU:H	1.60	0.66
1:A:14:ILE:HG23	1:A:15:GLN:H	1.60	0.66
2:O:114:ILE:O	2:O:116:ASN:N	2.28	0.66
2:O:18:LEU:HB2	2:O:31:ALA:HB2	1.77	0.66
4:Q:65:LYS:HD3	4:Q:68:LYS:HG2	1.78	0.66
1:A:26:VAL:CG2	2:B:29:GLU:HG3	2.25	0.66
4:D:66:VAL:HG12	4:D:67:SER:N	2.11	0.66
2:B:77:PRO:HG3	2:B:85:PHE:HB2	1.76	0.66
4:Q:150:LEU:N	4:Q:178:TRP:CZ2	2.63	0.66
11:A:305:PL9:H203	12:D:306:OPC:CBE	2.20	0.66
3:C:161:TYR:HE1	3:C:167:SER:CA	2.09	0.66
4:D:108:CYS:C	4:D:110:HIS:H	1.99	0.66
4:Q:150:LEU:N	4:Q:178:TRP:HZ2	1.94	0.66
3:C:28:ALA:O	3:C:239:GLY:HA3	1.94	0.66
1:A:158:ILE:CG2	1:A:159:PRO:HD2	2.25	0.66
4:D:105:ASN:O	4:D:107:VAL:N	2.29	0.66
3:C:70:LEU:H	3:C:70:LEU:HD23	1.61	0.66
4:Q:66:VAL:HG22	4:Q:160:ILE:HG13	1.76	0.66
3:P:1:TYR:HD2	3:P:119:LEU:HD21	1.59	0.66
3:P:120:PRO:HB2	3:P:124:TYR:HB2	1.77	0.66
4:D:73:HIS:NE2	4:D:91:ILE:HG22	2.11	0.65
4:Q:149:ALA:C	4:Q:150:LEU:HD22	2.17	0.65
3:P:180:ILE:HD11	3:P:182:LYS:O	1.96	0.65
4:Q:104:ILE:HA	4:Q:150:LEU:HD13	1.78	0.65
1:A:211:ILE:CD1	1:A:211:ILE:H	2.08	0.65
2:B:21:GLY:O	2:B:22:MET:HB2	1.97	0.65
3:C:177:THR:HA	3:C:227:ALA:HA	1.78	0.65
4:D:124:PHE:O	4:D:133:TYR:N	2.29	0.65
4:Q:78:ARG:HH12	4:Q:100:ARG:HH21	1.43	0.65
11:Q:1305:PL9:C21	11:Q:1305:PL9:H262	2.12	0.65
13:O:1201:CLA:HHC	13:O:1201:CLA:HBB1	1.76	0.65
2:B:80:TYR:HB2	13:B:201:CLA:HED1	1.77	0.65
3:C:120:PRO:HB2	3:C:124:TYR:HB2	1.79	0.65
1:N:114:ARG:HH21	1:N:209:GLN:H	1.41	0.65
1:N:27:PRO:HG3	2:O:33:PRO:CG	2.18	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:209:GLN:CG	1:A:210:GLY:N	2.53	0.65
2:B:118:ASN:HD21	2:B:126:ARG:HD2	1.60	0.65
3:P:188:ASP:HB3	3:P:192:ASN:O	1.95	0.65
6:S:20:TRP:O	6:S:24:VAL:HG23	1.96	0.65
4:Q:168:THR:HG23	4:Q:174:GLU:C	2.17	0.65
1:A:27:PRO:HG3	2:B:33:PRO:CD	2.27	0.65
4:D:155:VAL:HA	4:D:160:ILE:HA	1.79	0.65
3:P:90:ILE:CG2	3:P:95:LYS:HB2	2.26	0.65
1:A:106:LEU:HD12	5:E:18:ILE:CD1	2.27	0.65
3:P:218:ILE:HG13	3:P:219:VAL:H	1.60	0.65
6:F:25:LEU:O	6:F:25:LEU:HD23	1.96	0.65
2:B:76:LEU:H	2:B:76:LEU:HD22	1.62	0.65
1:A:26:VAL:HB	2:B:29:GLU:HG3	1.79	0.65
3:C:281:LYS:HA	3:C:284:ALA:CB	2.19	0.65
3:C:188:ASP:HB3	3:C:192:ASN:O	1.97	0.65
4:D:134:ASP:HB3	4:D:138:ARG:CG	2.27	0.65
5:E:5:ALA:CB	6:F:5:MET:HB2	2.27	0.65
3:C:172:PHE:N	3:C:231:LEU:HG	2.02	0.65
3:P:277:LYS:HE3	7:T:31:ARG:O	1.96	0.65
6:S:26:LEU:CD2	8:U:17:TRP:HE1	2.09	0.65
2:O:74:GLU:HG2	2:O:75:ILE:H	1.61	0.65
11:Q:1305:PL9:H302	12:Q:1307:OPC:CBC	2.27	0.65
3:C:231:LEU:HD11	3:C:233:ASN:O	1.96	0.65
4:D:66:VAL:HG23	4:D:158:ASP:C	2.17	0.65
1:A:36:LEU:HD23	1:A:100:HIS:HA	1.79	0.65
10:A:304:TDS:CBB	2:B:88:LEU:HD22	2.26	0.65
3:P:120:PRO:HB3	3:P:124:TYR:CD1	2.32	0.65
1:N:103:ARG:HG3	1:N:103:ARG:HH11	1.61	0.65
5:R:16:PHE:HB3	15:R:1101:BCR:C36	2.20	0.65
8:H:19:ILE:O	8:H:23:VAL:HG13	1.97	0.65
2:O:146:GLY:O	2:O:149:LEU:HG	1.97	0.65
1:A:157:ALA:HB3	2:B:98:VAL:HG21	1.79	0.65
3:P:226:LYS:HG2	3:P:229:GLU:HB3	1.79	0.65
4:D:81:VAL:HG12	4:D:82:GLN:N	2.11	0.64
4:D:65:LYS:HZ2	4:D:158:ASP:HB3	1.61	0.64
3:P:185:LYS:HE2	3:P:195:TYR:CB	2.28	0.64
2:O:32:TRP:HZ3	2:O:36:LEU:HA	1.62	0.64
4:Q:14:GLY:O	4:Q:16:ARG:HG2	1.96	0.64
7:G:14:VAL:O	7:G:17:THR:HB	1.95	0.64
7:T:26:TYR:H	7:T:26:TYR:HD1	1.44	0.64
3:P:281:LYS:CA	3:P:284:ALA:HB3	2.21	0.64
1:N:47:GLN:HE22	1:N:89:SER:CB	2.11	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:D:14:GLY:O	4:D:16:ARG:HG2	1.96	0.64
2:O:27:TYR:HD2	2:O:27:TYR:O	1.80	0.64
4:Q:125:LYS:HD3	4:Q:132:GLN:HG2	1.80	0.64
4:Q:76:GLY:H	4:Q:93:VAL:CG1	2.01	0.64
1:A:24:LYS:HD3	1:A:24:LYS:N	2.11	0.64
1:A:156:GLU:HA	1:A:162:GLY:O	1.97	0.64
3:P:177:THR:HA	3:P:227:ALA:HA	1.78	0.64
4:D:73:HIS:HE2	4:D:91:ILE:HG22	1.62	0.64
3:C:273:ILE:HD13	7:G:30:LYS:CD	2.14	0.64
3:C:273:ILE:CD1	7:G:30:LYS:HD3	2.12	0.64
3:P:54:TYR:HD2	3:P:70:LEU:HD13	1.63	0.64
4:Q:115:VAL:HG22	4:Q:126:CYS:HB2	1.79	0.64
4:Q:153:ALA:N	4:Q:162:LEU:HG	2.12	0.64
2:B:26:TYR:H	2:B:26:TYR:HD2	1.45	0.64
2:O:122:ASN:CG	5:R:30:LYS:HD3	2.17	0.64
4:Q:133:TYR:HA	4:Q:139:VAL:CA	2.27	0.64
3:C:218:ILE:HG12	3:C:232:THR:HA	1.80	0.64
4:D:66:VAL:HB	4:D:158:ASP:HA	1.80	0.64
1:A:136:TYR:CD1	2:B:78:GLU:HG2	2.33	0.64
4:D:121:GLU:OE2	4:D:125:LYS:HG3	1.98	0.64
6:F:22:LEU:O	6:F:26:LEU:HD23	1.98	0.64
6:F:27:LEU:O	6:F:30:GLN:HG3	1.96	0.64
8:H:4:ASP:O	8:H:8:TRP:CB	2.45	0.64
3:C:216:GLU:O	3:C:233:ASN:HB2	1.96	0.64
1:A:142:GLN:NE2	2:B:67:ASN:N	2.45	0.64
3:P:107:LYS:HD3	3:P:110:GLN:NE2	2.12	0.64
3:P:126:GLU:HA	3:P:126:GLU:OE1	1.96	0.64
3:P:79:PRO:HG2	3:P:82:PHE:CE1	2.33	0.64
1:N:111:LYS:CB	1:N:114:ARG:HH11	2.11	0.64
1:N:28:PRO:HG2	2:O:32:TRP:HB2	1.80	0.64
1:N:31:ASN:HB3	1:N:34:TYR:CG	2.32	0.64
4:Q:64:VAL:HG12	4:Q:91:ILE:HG13	1.78	0.64
4:D:105:ASN:O	4:D:107:VAL:HG23	1.98	0.64
3:C:107:LYS:HD3	3:C:110:GLN:NE2	2.13	0.64
4:Q:15:ARG:NH2	4:Q:17:GLN:HG2	2.12	0.64
4:D:136:THR:CB	4:D:138:ARG:HD2	2.16	0.64
4:Q:94:GLU:HG2	4:Q:95:SER:H	1.62	0.64
3:P:218:ILE:HG12	3:P:232:THR:HA	1.79	0.64
2:O:22:MET:C	2:O:24:HIS:N	2.50	0.64
2:B:26:TYR:C	2:B:27:TYR:CD1	2.72	0.64
3:C:57:LYS:O	3:C:58:LEU:HB2	1.98	0.64
2:B:91:LEU:HD22	2:B:96:LEU:HD13	1.80	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:Q:134:ASP:HB2	4:Q:138:ARG:HB2	1.80	0.63
2:O:57:LEU:HD21	3:P:258:MET:CE	2.28	0.63
3:P:231:LEU:CD1	3:P:232:THR:H	2.11	0.63
2:O:24:HIS:O	2:O:25:ASN:ND2	2.31	0.63
3:P:57:LYS:O	3:P:58:LEU:HB2	1.97	0.63
1:A:146:TRP:HB2	2:B:75:ILE:HD11	1.79	0.63
2:B:122:ASN:ND2	2:B:124:PHE:HD2	1.96	0.63
8:H:10:ALA:O	8:H:14:VAL:HG23	1.97	0.63
1:A:177:GLN:HB2	1:N:59:LYS:HZ3	1.61	0.63
3:C:81:GLY:C	3:C:134:PRO:HG3	2.19	0.63
4:Q:176:PRO:HB2	4:Q:178:TRP:CZ3	2.33	0.63
1:A:27:PRO:HG3	2:B:33:PRO:HD3	1.80	0.63
3:C:185:LYS:HD3	3:C:195:TYR:HD2	1.64	0.63
3:C:234:ASN:OD1	3:C:236:ASN:HB3	1.97	0.63
3:P:171:VAL:CG1	3:P:234:ASN:HB2	2.27	0.63
2:O:20:LYS:CG	2:O:21:GLY:N	2.61	0.63
4:D:132:GLN:N	4:D:144:ALA:HB2	2.13	0.63
2:B:122:ASN:HD21	2:B:124:PHE:HB3	1.60	0.63
3:C:270:LEU:HA	7:G:27:GLN:OE1	1.98	0.63
4:D:66:VAL:HG22	4:D:160:ILE:CG1	2.03	0.63
3:P:265:VAL:O	3:P:269:GLN:HG3	1.98	0.63
3:C:218:ILE:HG13	3:C:232:THR:HG22	1.81	0.63
3:P:38:GLN:OE1	7:T:15:PHE:HE1	1.80	0.63
8:U:19:ILE:O	8:U:23:VAL:HG13	1.99	0.63
6:S:27:LEU:O	6:S:30:GLN:HG3	1.99	0.63
2:O:124:PHE:CE2	5:R:26:ILE:HB	2.33	0.63
4:Q:138:ARG:NH2	4:Q:147:SER:OG	2.31	0.63
6:F:26:LEU:HD21	8:H:17:TRP:HE1	1.62	0.63
8:H:23:VAL:HA	8:H:26:ARG:HD2	1.81	0.63
2:O:133:PHE:CD2	13:O:1201:CLA:HMB3	2.32	0.63
3:C:89:ARG:O	3:C:91:PRO:HD3	1.98	0.63
4:D:134:ASP:OD1	4:D:135:GLU:N	2.32	0.63
4:D:150:LEU:CD1	4:D:171:ARG:HD3	2.28	0.63
4:D:166:THR:CA	4:D:179:VAL:HG13	2.13	0.63
4:Q:121:GLU:OE1	4:Q:125:LYS:HE3	1.99	0.63
4:Q:24:PHE:CD1	11:Q:1305:PL9:H202	2.34	0.63
2:O:114:ILE:C	2:O:116:ASN:H	2.02	0.63
4:Q:165:TRP:CZ2	4:Q:176:PRO:HB3	2.33	0.63
3:C:262:ILE:HG21	7:G:17:THR:CG2	2.29	0.63
7:G:25:ALA:HB1	7:G:29:TYR:OH	1.99	0.63
2:O:67:ASN:HD21	3:P:16:PRO:HB3	1.63	0.63
1:A:177:GLN:HB2	1:N:59:LYS:NZ	2.13	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
6:F:26:LEU:HB3	7:G:30:LYS:HE3	1.81	0.62
5:R:3:LEU:HA	5:R:6:VAL:CG2	2.29	0.62
2:O:117:VAL:O	2:O:118:ASN:CB	2.47	0.62
4:Q:168:THR:HA	4:Q:176:PRO:CD	2.13	0.62
11:A:305:PL9:H462	11:A:305:PL9:C41	2.12	0.62
2:O:146:GLY:C	2:O:149:LEU:HG	2.19	0.62
7:T:26:TYR:CD1	7:T:26:TYR:N	2.65	0.62
1:N:13:GLU:OE1	1:N:14:ILE:HG22	1.98	0.62
4:Q:155:VAL:HG12	4:Q:156:GLN:N	2.14	0.62
1:A:26:VAL:HG21	2:B:29:GLU:HG3	1.81	0.62
5:R:3:LEU:HA	5:R:6:VAL:HG23	1.81	0.62
3:C:226:LYS:HE3	3:C:229:GLU:C	2.19	0.62
1:A:62:VAL:HG13	1:A:177:GLN:HB3	1.79	0.62
1:N:27:PRO:O	2:O:29:GLU:HB3	1.99	0.62
4:Q:116:PRO:CD	4:Q:126:CYS:HA	2.24	0.62
3:C:271:MET:CG	4:D:22:LEU:HD21	2.23	0.62
1:N:45:LEU:HB3	11:Q:1305:PL9:H502	1.81	0.62
3:C:93:GLU:CA	3:C:97:GLU:HB2	2.23	0.62
3:P:94:LEU:CD2	3:P:94:LEU:H	2.06	0.62
1:A:118:TRP:HD1	9:A:302:HEM:HMD2	1.64	0.62
4:D:51:GLY:CA	4:D:164:PRO:HG2	2.21	0.62
3:P:83:LYS:HB2	3:P:111:ASP:CB	2.29	0.62
4:D:117:TRP:HD1	4:D:124:PHE:CE1	2.17	0.62
12:B:307:OPC:HBU2	12:B:307:OPC:HBC1	1.80	0.62
5:E:23:ILE:HG23	5:E:24:PHE:N	2.14	0.62
12:Q:1307:OPC:HBZ2	12:Q:1307:OPC:HBF1	1.80	0.62
3:P:10:PRO:HG2	3:P:11:PRO:HD3	1.80	0.62
4:Q:19:MET:O	4:Q:22:LEU:HB3	1.99	0.62
1:A:26:VAL:CB	2:B:29:GLU:HG3	2.29	0.62
2:B:32:TRP:CE3	2:B:36:LEU:HG	2.35	0.62
3:P:89:ARG:O	3:P:91:PRO:HD3	2.00	0.62
3:C:10:PRO:HG2	3:C:11:PRO:HD3	1.82	0.62
5:R:5:ALA:CB	6:S:5:MET:HB2	2.30	0.62
3:C:185:LYS:HE2	3:C:195:TYR:HB3	1.81	0.62
4:D:121:GLU:OE1	4:D:125:LYS:HE3	2.00	0.62
5:E:27:LYS:O	5:E:31:LEU:HD22	1.99	0.62
12:Q:1307:OPC:HAL2	12:Q:1307:OPC:HAP1	1.82	0.62
3:C:218:ILE:HG13	3:C:219:VAL:N	2.14	0.62
2:B:149:LEU:HB2	2:B:150:PRO:CD	2.29	0.62
2:O:20:LYS:CD	2:O:21:GLY:H	2.13	0.62
5:E:11:PHE:C	5:E:11:PHE:CD1	2.72	0.62
2:O:122:ASN:HD21	5:R:30:LYS:HG3	1.64	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:D:90:TYR:CE2	4:D:106:ALA:HB2	2.34	0.62
4:Q:134:ASP:CB	4:Q:138:ARG:HB2	2.30	0.62
7:G:26:TYR:CD1	7:G:26:TYR:N	2.67	0.62
7:T:28:GLN:O	7:T:30:LYS:N	2.33	0.62
7:T:9:LEU:HD23	7:T:9:LEU:N	2.15	0.62
3:P:111:ASP:O	3:P:114:LEU:HD21	2.00	0.62
1:N:21:VAL:HG12	1:N:22:THR:N	2.15	0.62
1:N:21:VAL:O	1:N:23:SER:N	2.32	0.62
3:P:275:LYS:CD	4:Q:20:ASN:HB3	2.30	0.61
5:R:23:ILE:HG23	5:R:24:PHE:N	2.14	0.61
2:B:18:LEU:O	2:B:19:ALA:HB3	1.99	0.61
4:D:121:GLU:HB3	4:D:123:LYS:HG3	1.80	0.61
4:D:73:HIS:ND1	4:D:77:ASP:HB3	2.15	0.61
5:R:11:PHE:C	5:R:11:PHE:CD1	2.73	0.61
4:D:111:LEU:HB3	4:D:129:HIS:CE1	2.36	0.61
2:O:144:GLY:O	2:O:145:ILE:HD13	2.00	0.61
12:N:1306:OPC:HAP2	12:N:1306:OPC:CBI	2.23	0.61
3:P:188:ASP:HB3	3:P:193:VAL:HA	1.82	0.61
2:O:20:LYS:CG	2:O:21:GLY:H	2.13	0.61
4:D:136:THR:HG22	4:D:150:LEU:HD11	1.83	0.61
1:A:30:VAL:HG21	5:E:26:ILE:HD11	1.82	0.61
3:P:117:GLY:O	3:P:119:LEU:HG	2.01	0.61
1:A:15:GLN:O	1:A:16:ALA:HB3	2.00	0.61
3:C:171:VAL:HB	3:C:231:LEU:CD1	2.30	0.61
3:C:226:LYS:HG2	3:C:229:GLU:CB	2.30	0.61
3:P:81:GLY:C	3:P:134:PRO:HG3	2.20	0.61
1:A:116:LEU:N	1:A:116:LEU:HD12	2.14	0.61
10:N:1304:TDS:HAU1	13:O:1201:CLA:H91	1.81	0.61
7:T:26:TYR:N	7:T:26:TYR:HD1	1.97	0.61
1:A:105:TYR:HD2	1:A:106:LEU:HD23	1.65	0.61
2:O:134:LEU:O	2:O:138:LEU:HD13	2.01	0.61
5:R:21:GLY:O	5:R:25:ALA:HB2	2.00	0.61
3:C:188:ASP:O	3:C:189:GLU:HB2	2.01	0.61
3:C:188:ASP:HB3	3:C:193:VAL:HA	1.81	0.61
3:P:231:LEU:HD13	3:P:232:THR:N	2.14	0.61
3:P:226:LYS:HE3	3:P:229:GLU:C	2.20	0.61
9:N:301:HEM:CHA	9:N:301:HEM:HBD1	2.31	0.61
1:N:158:ILE:O	1:N:162:GLY:HA3	2.01	0.61
4:Q:104:ILE:CD1	4:Q:136:THR:HA	2.30	0.61
4:D:65:LYS:HD3	4:D:68:LYS:NZ	2.16	0.61
12:N:1306:OPC:HBQ2	12:N:1306:OPC:HAT1	1.82	0.61
2:O:124:PHE:HZ	5:R:26:ILE:CB	2.10	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:N:98:ILE:HD11	13:O:1201:CLA:HED3	1.81	0.61
1:N:196:ALA:O	1:N:200:LEU:HB2	2.00	0.61
3:P:226:LYS:HE3	3:P:229:GLU:HB3	1.81	0.61
2:B:75:ILE:O	2:B:76:LEU:O	2.18	0.60
4:Q:123:LYS:HE3	4:Q:125:LYS:HZ3	1.64	0.60
4:Q:172:THR:HB	4:Q:174:GLU:OE2	2.01	0.60
4:Q:65:LYS:HB2	4:Q:68:LYS:HB2	1.82	0.60
1:A:200:LEU:O	1:A:204:LEU:HG	2.00	0.60
3:P:91:PRO:HD2	3:P:95:LYS:CG	2.31	0.60
5:E:18:ILE:O	5:E:22:ILE:HB	2.01	0.60
1:N:116:LEU:N	1:N:116:LEU:HD12	2.14	0.60
2:O:118:ASN:ND2	2:O:121:GLN:HE22	1.95	0.60
1:N:214:PRO:CB	5:R:29:ILE:HG22	2.13	0.60
3:C:171:VAL:HG12	3:C:233:ASN:O	2.01	0.60
2:B:112:PRO:O	2:B:116:ASN:CB	2.49	0.60
1:N:110:PHE:HE1	2:O:111:VAL:HG12	1.67	0.60
4:Q:58:ASP:OD1	4:Q:64:VAL:HG22	2.01	0.60
4:Q:56:ALA:CA	4:Q:81:VAL:HG13	2.21	0.60
3:C:86:PRO:HG2	3:C:89:ARG:HG3	1.83	0.60
2:O:75:ILE:O	2:O:77:PRO:HD2	2.01	0.60
1:N:37:GLY:HA3	9:N:302:HEM:HBA1	1.82	0.60
4:D:136:THR:HB	4:D:138:ARG:CD	2.17	0.60
4:D:168:THR:CA	4:D:176:PRO:HD3	2.31	0.60
4:D:42:PHE:CD1	12:D:306:OPC:HAG2	2.36	0.60
5:R:5:ALA:HA	5:R:8:TYR:HB2	1.84	0.60
3:C:275:LYS:CE	4:D:17:GLN:HB2	2.32	0.60
3:C:90:ILE:CG2	3:C:95:LYS:HB2	2.32	0.60
3:C:83:LYS:HB2	3:C:111:ASP:CB	2.30	0.60
3:P:45:VAL:HG22	3:P:85:ALA:HB2	1.83	0.60
3:P:118:PRO:C	3:P:120:PRO:HD3	2.22	0.60
3:P:275:LYS:HD3	3:P:275:LYS:C	2.21	0.60
4:D:123:LYS:HD2	4:D:140:ILE:HD11	1.83	0.60
4:D:167:GLU:C	4:D:176:PRO:HG3	2.21	0.60
4:Q:81:VAL:HB	4:Q:89:THR:OG1	2.01	0.60
4:D:19:MET:O	4:D:22:LEU:HB3	2.01	0.60
2:O:106:LEU:HD11	13:O:1201:CLA:H171	1.84	0.60
3:C:180:ILE:O	3:C:221:GLU:HA	2.02	0.60
1:N:28:PRO:HG2	2:O:32:TRP:CB	2.31	0.60
3:C:262:ILE:HG21	7:G:17:THR:HG23	1.82	0.60
11:A:305:PL9:H252	12:D:306:OPC:CBN	2.31	0.60
2:O:141:ILE:HD11	5:R:11:PHE:CZ	2.36	0.60
4:D:51:GLY:C	4:D:85:LYS:HZ1	2.04	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:175:VAL:O	1:A:175:VAL:HG12	2.01	0.60
15:R:1101:BCR:H362	6:S:17:PHE:HZ	1.67	0.60
2:O:122:ASN:CG	5:R:26:ILE:HG21	2.21	0.60
1:A:204:LEU:C	1:A:206:ILE:H	2.04	0.60
6:F:11:LEU:O	6:F:15:LEU:HD23	2.02	0.60
3:C:218:ILE:HG12	3:C:232:THR:CA	2.31	0.60
3:P:180:ILE:O	3:P:221:GLU:HA	2.02	0.60
3:P:218:ILE:HG13	3:P:219:VAL:N	2.17	0.60
4:D:100:ARG:O	4:D:153:ALA:HB2	2.02	0.60
4:D:101:ASP:HA	4:D:153:ALA:HB2	1.82	0.60
4:Q:129:HIS:HB2	14:Q:1200:FES:S1	2.42	0.60
3:C:116:VAL:O	3:C:116:VAL:HG12	2.00	0.60
4:D:137:GLY:CA	4:D:148:LEU:HB2	2.32	0.59
2:B:124:PHE:HE1	5:E:23:ILE:HG13	1.67	0.59
2:B:38:TYR:OH	12:B:307:OPC:HAS2	2.01	0.59
2:B:42:VAL:HG23	3:C:272:LEU:HD13	1.82	0.59
1:N:191:LEU:HD21	10:N:1304:TDS:OBD	2.02	0.59
4:D:66:VAL:HG11	4:D:155:VAL:HG11	1.83	0.59
3:C:55:ASP:OD1	3:C:57:LYS:HG3	2.02	0.59
4:D:13:MET:N	4:D:13:MET:SD	2.75	0.59
3:C:266:MET:HE3	7:G:20:GLY:O	2.02	0.59
7:G:26:TYR:CA	7:G:29:TYR:CD1	2.82	0.59
2:O:146:GLY:O	2:O:149:LEU:N	2.31	0.59
2:O:126:ARG:HH12	2:O:128:VAL:C	2.06	0.59
2:B:70:ALA:HB1	3:C:17:THR:HA	1.83	0.59
3:P:283:GLN:C	3:P:285:ALA:H	2.04	0.59
1:N:39:ILE:HD11	15:R:1101:BCR:H312	1.83	0.59
4:D:121:GLU:O	4:D:122:ASN:HB2	2.03	0.59
4:Q:66:VAL:O	4:Q:70:LEU:HB2	2.02	0.59
12:D:306:OPC:HAY2	12:D:306:OPC:HBC2	1.83	0.59
1:A:95:LEU:HD13	1:A:96:MET:CE	2.32	0.59
3:P:217:LEU:CA	3:P:232:THR:HB	2.32	0.59
4:Q:96:LYS:O	4:Q:97:GLU:HB2	2.01	0.59
4:D:138:ARG:NE	4:D:171:ARG:HD2	2.18	0.59
11:A:305:PL9:H111	12:D:306:OPC:CBF	2.32	0.59
2:B:34:ASN:ND2	12:B:307:OPC:CBK	2.64	0.59
7:G:11:LEU:HA	7:G:14:VAL:HG11	1.84	0.59
3:P:194:LYS:HE3	3:P:212:PRO:HB3	1.83	0.59
3:C:45:VAL:HG22	3:C:85:ALA:HB2	1.84	0.59
4:D:132:GLN:H	4:D:144:ALA:HB2	1.68	0.59
4:Q:94:GLU:HG2	4:Q:95:SER:N	2.18	0.59
7:G:11:LEU:HA	7:G:14:VAL:HG13	1.85	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:P:278:GLN:O	3:P:281:LYS:HE2	2.03	0.59
3:P:28:ALA:O	3:P:239:GLY:HA3	2.03	0.59
3:P:105:PRO:HB3	3:P:110:GLN:O	2.02	0.59
2:B:71:THR:CB	2:B:72:PRO:HD3	2.11	0.59
4:Q:66:VAL:HG22	4:Q:160:ILE:HG12	1.85	0.59
3:C:171:VAL:HG13	3:C:234:ASN:HD22	1.68	0.59
3:P:40:VAL:HG13	3:P:247:ILE:HD11	1.85	0.59
4:D:123:LYS:HZ1	4:D:140:ILE:HG13	1.67	0.59
4:Q:91:ILE:HG23	4:Q:160:ILE:CD1	2.32	0.59
4:Q:55:THR:O	4:Q:56:ALA:HB3	2.03	0.59
2:B:32:TRP:CZ3	2:B:36:LEU:HG	2.37	0.59
3:C:269:GLN:NE2	7:G:24:ALA:HA	2.17	0.59
5:E:3:LEU:HA	5:E:6:VAL:HG23	1.85	0.59
6:F:12:SER:O	6:F:16:ILE:HG12	2.02	0.59
1:N:204:LEU:C	1:N:206:ILE:H	2.05	0.59
3:P:195:TYR:O	3:P:196:GLN:O	2.21	0.59
2:B:86:GLN:HG2	2:B:143:LEU:HA	1.84	0.59
4:D:99:ILE:O	4:D:99:ILE:HD12	2.03	0.59
4:Q:93:VAL:HG22	4:Q:94:GLU:N	2.17	0.59
8:H:17:TRP:O	8:H:20:ALA:HB3	2.03	0.59
4:D:65:LYS:O	4:D:68:LYS:HB2	2.03	0.59
7:T:31:ARG:HD3	7:T:32:PRO:HD3	1.85	0.59
3:P:86:PRO:HG2	3:P:89:ARG:HG3	1.84	0.59
3:P:216:GLU:HB2	3:P:233:ASN:OD1	2.03	0.59
3:C:54:TYR:HD2	3:C:70:LEU:HD13	1.66	0.59
1:N:81:LEU:HD13	1:N:81:LEU:C	2.23	0.59
3:C:49:VAL:HG22	3:C:128:VAL:HG22	1.85	0.59
5:R:16:PHE:O	5:R:20:VAL:HB	2.03	0.59
4:Q:64:VAL:HG21	4:Q:80:LEU:C	2.23	0.59
8:H:11:LEU:HB3	8:H:15:PHE:HZ	1.66	0.59
5:E:5:ALA:HA	5:E:8:TYR:HB2	1.85	0.59
1:N:154:VAL:HG11	10:N:1304:TDS:HAX1	1.84	0.59
10:A:304:TDS:CBC	2:B:85:PHE:HD2	2.15	0.59
2:O:25:ASN:OD1	2:O:26:TYR:N	2.35	0.59
5:R:9:ILE:HG12	6:S:13:PHE:HB2	1.85	0.59
4:D:166:THR:C	4:D:179:VAL:HG22	2.23	0.59
4:Q:134:ASP:OD1	4:Q:138:ARG:HB2	2.03	0.59
4:Q:138:ARG:HB3	4:Q:138:ARG:NH1	2.18	0.59
4:Q:156:GLN:CB	4:Q:161:VAL:HG21	2.23	0.59
1:A:35:CYS:O	1:A:39:ILE:HG12	2.02	0.59
5:E:26:ILE:HA	5:E:29:ILE:HG13	1.84	0.59
7:T:17:THR:O	7:T:21:LEU:HB2	2.03	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:P:178:GLY:C	3:P:225:VAL:HA	2.23	0.58
4:D:131:SER:OG	4:D:144:ALA:HA	2.02	0.58
2:B:118:ASN:CB	2:B:121:GLN:O	2.49	0.58
3:C:194:LYS:HE3	3:C:212:PRO:HB3	1.85	0.58
2:B:149:LEU:N	2:B:149:LEU:HD12	2.18	0.58
4:D:80:LEU:HD13	4:D:88:PRO:HB2	1.85	0.58
3:P:116:VAL:HG12	3:P:116:VAL:O	2.03	0.58
3:P:49:VAL:HG22	3:P:128:VAL:HG22	1.85	0.58
4:Q:66:VAL:HG12	4:Q:70:LEU:HD12	1.85	0.58
1:A:23:SER:HB2	1:A:24:LYS:HD3	1.86	0.58
12:Q:1307:OPC:CAX	12:Q:1307:OPC:HAT1	2.33	0.58
7:T:27:GLN:HA	7:T:27:GLN:NE2	2.18	0.58
3:C:59:GLN:CB	3:C:67:LYS:HB3	2.33	0.58
3:P:275:LYS:CE	4:Q:17:GLN:HB2	2.32	0.58
15:R:1101:BCR:H362	6:S:17:PHE:CZ	2.38	0.58
4:D:134:ASP:N	4:D:138:ARG:O	2.36	0.58
4:D:104:ILE:CG2	4:D:148:LEU:HB3	2.34	0.58
4:Q:142:GLY:CA	4:Q:144:ALA:H	2.17	0.58
11:A:305:PL9:H162	12:B:307:OPC:HBU1	1.84	0.58
7:G:30:LYS:O	7:G:31:ARG:CZ	2.50	0.58
1:A:33:PHE:CE1	5:E:14:LEU:HD22	2.39	0.58
3:P:218:ILE:HG12	3:P:232:THR:CA	2.33	0.58
3:C:226:LYS:CD	3:C:229:GLU:HB3	2.33	0.58
2:O:24:HIS:N	2:O:24:HIS:ND1	2.50	0.58
1:N:23:SER:HA	1:N:25:TYR:CE2	2.39	0.58
3:C:50:VAL:HG13	3:C:151:LEU:HD21	1.85	0.58
4:Q:78:ARG:NH2	4:Q:100:ARG:NH2	2.46	0.58
4:Q:133:TYR:HB3	4:Q:137:GLY:C	2.23	0.58
1:N:95:LEU:HD13	1:N:96:MET:CE	2.33	0.58
3:C:216:GLU:O	3:C:233:ASN:CB	2.51	0.58
3:C:91:PRO:HD2	3:C:95:LYS:CG	2.33	0.58
4:D:50:VAL:HG12	4:D:164:PRO:HB3	1.84	0.58
3:P:231:LEU:HD12	3:P:231:LEU:H	1.68	0.58
4:D:136:THR:O	4:D:136:THR:HG22	2.04	0.58
4:Q:168:THR:HG23	4:Q:175:LYS:N	2.17	0.58
1:A:196:ALA:O	1:A:200:LEU:HB2	2.04	0.58
3:C:255:VAL:CG1	7:G:14:VAL:CG1	2.74	0.58
8:H:12:LEU:HA	8:H:15:PHE:CD1	2.38	0.58
3:P:59:GLN:CB	3:P:67:LYS:HB3	2.34	0.58
6:F:7:TYR:O	6:F:11:LEU:HD12	2.04	0.58
7:G:26:TYR:HD1	7:G:26:TYR:N	2.00	0.58
8:U:7:GLY:O	8:U:11:LEU:HD11	2.03	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
8:U:23:VAL:O	8:U:27:ASN:ND2	2.36	0.58
4:D:137:GLY:HA3	4:D:148:LEU:HB2	1.85	0.58
4:D:129:HIS:O	4:D:143:PRO:HG3	2.03	0.58
7:G:9:LEU:HG	7:G:10:VAL:H	1.69	0.58
3:C:120:PRO:HB3	3:C:124:TYR:CD1	2.39	0.58
3:C:261:PHE:CE2	4:D:30:VAL:HG13	2.38	0.58
4:Q:152:HIS:H	4:Q:162:LEU:HD23	1.69	0.58
8:U:6:LEU:C	8:U:8:TRP:N	2.57	0.58
1:A:193:TRP:O	1:A:197:VAL:HG23	2.03	0.58
4:D:116:PRO:CG	4:D:126:CYS:HA	2.34	0.58
4:D:93:VAL:HG13	4:D:94:GLU:O	2.03	0.58
2:B:38:TYR:HB3	3:C:276:LYS:CG	2.32	0.58
7:T:21:LEU:C	7:T:23:TYR:H	2.05	0.58
3:P:188:ASP:O	3:P:189:GLU:HB2	2.03	0.58
1:N:188:THR:CG2	9:N:301:HEM:HBC2	2.33	0.58
2:O:118:ASN:HD22	2:O:121:GLN:CD	2.06	0.57
2:O:122:ASN:CG	5:R:26:ILE:CG2	2.72	0.57
11:A:305:PL9:H513	4:D:28:THR:HG23	1.85	0.57
7:G:26:TYR:HA	7:G:29:TYR:HD1	1.61	0.57
8:H:6:LEU:C	8:H:8:TRP:H	2.05	0.57
4:D:134:ASP:HB3	4:D:138:ARG:HG2	1.85	0.57
4:Q:123:LYS:HB2	4:Q:125:LYS:HE2	1.86	0.57
12:Q:1307:OPC:CAQ	12:Q:1307:OPC:HBG3	2.34	0.57
8:U:6:LEU:HA	8:U:9:VAL:HG12	1.86	0.57
2:B:119:LYS:O	2:B:119:LYS:HD3	2.03	0.57
4:Q:178:TRP:N	4:Q:178:TRP:CE3	2.69	0.57
2:O:84:VAL:HG22	13:O:1201:CLA:HMD1	1.86	0.57
3:C:3:PHE:HZ	3:C:119:LEU:HD11	1.69	0.57
1:N:112:LYS:HB2	1:N:113:PRO:CD	2.27	0.57
1:N:113:PRO:C	1:N:115:GLU:OE1	2.42	0.57
1:N:29:HIS:HE2	1:N:209:GLN:CG	2.16	0.57
4:Q:88:PRO:HG3	4:Q:106:ALA:N	2.20	0.57
4:Q:123:LYS:O	4:Q:125:LYS:HG2	2.03	0.57
7:G:25:ALA:O	7:G:29:TYR:CD2	2.57	0.57
8:H:6:LEU:C	8:H:8:TRP:N	2.56	0.57
3:P:255:VAL:CG1	7:T:14:VAL:HG12	2.34	0.57
3:P:52:ILE:HG12	3:P:153:ALA:CB	2.34	0.57
4:D:84:LEU:O	4:D:86:GLY:N	2.37	0.57
4:D:150:LEU:HD13	4:D:171:ARG:HD3	1.86	0.57
1:A:27:PRO:HG3	2:B:33:PRO:HG3	1.87	0.57
3:P:217:LEU:HA	3:P:232:THR:CB	2.33	0.57
3:P:84:ILE:HD12	3:P:130:PRO:HD2	1.86	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:D:110:HIS:CE1	4:D:129:HIS:HB2	2.39	0.57
4:Q:64:VAL:O	4:Q:91:ILE:HD11	2.04	0.57
3:C:278:GLN:O	3:C:281:LYS:HE2	2.04	0.57
3:P:202:ASP:HA	3:P:206:THR:CB	2.25	0.57
1:A:13:GLU:C	1:A:17:LEU:HD12	2.25	0.57
4:D:124:PHE:CB	4:D:133:TYR:HB2	2.22	0.57
4:D:65:LYS:HD3	4:D:68:LYS:HZ1	1.69	0.57
3:P:258:MET:O	3:P:262:ILE:HD13	2.04	0.57
3:P:216:GLU:O	3:P:233:ASN:HB2	2.03	0.57
3:C:117:GLY:O	3:C:119:LEU:HG	2.05	0.57
2:O:72:PRO:O	2:O:73:LEU:HB3	2.04	0.57
3:C:79:PRO:HG2	3:C:82:PHE:CE1	2.39	0.57
3:C:275:LYS:HD3	3:C:275:LYS:C	2.25	0.57
3:C:231:LEU:HD22	3:C:232:THR:N	2.19	0.57
3:C:202:ASP:HA	3:C:206:THR:CB	2.27	0.57
3:P:61:VAL:O	3:P:157:ARG:NH1	2.37	0.57
1:A:62:VAL:HG13	1:A:177:GLN:CB	2.34	0.57
2:B:114:ILE:C	2:B:116:ASN:H	2.07	0.57
1:N:79:GLY:HA2	4:Q:41:TYR:HE1	1.69	0.57
4:D:128:CYS:HB3	4:D:129:HIS:CE1	2.40	0.57
4:D:166:THR:HG23	4:D:179:VAL:HG11	1.86	0.57
1:A:207:ARG:O	1:A:208:LYS:HD2	2.05	0.57
3:C:276:LYS:HE3	3:C:276:LYS:HA	1.87	0.57
3:P:91:PRO:HD2	3:P:95:LYS:HG2	1.86	0.57
3:P:172:PHE:C	3:P:231:LEU:HB3	2.25	0.57
2:O:74:GLU:O	2:O:75:ILE:HB	2.05	0.57
2:B:134:LEU:O	2:B:138:LEU:HD13	2.05	0.57
4:Q:116:PRO:HD2	4:Q:126:CYS:CA	2.26	0.57
2:B:32:TRP:H	2:B:33:PRO:HD3	1.70	0.57
3:C:271:MET:HB3	4:D:23:ALA:HA	1.87	0.57
5:E:27:LYS:O	5:E:31:LEU:HB2	2.05	0.57
8:H:23:VAL:O	8:H:27:ASN:ND2	2.37	0.57
5:E:3:LEU:HA	5:E:6:VAL:CG2	2.34	0.57
1:N:111:LYS:CB	1:N:114:ARG:NH1	2.68	0.56
1:N:33:PHE:HB3	1:N:103:ARG:HG2	1.87	0.56
1:N:209:GLN:CB	2:O:27:TYR:HB2	2.35	0.56
2:B:18:LEU:O	2:B:19:ALA:CB	2.51	0.56
2:O:80:TYR:CD1	2:O:80:TYR:N	2.73	0.56
1:N:95:LEU:HG	5:R:7:PHE:CB	2.34	0.56
2:B:84:VAL:O	2:B:88:LEU:HD13	2.04	0.56
3:C:269:GLN:CD	7:G:24:ALA:HA	2.26	0.56
11:Q:1305:PL9:C32	12:Q:1307:OPC:HBW2	2.34	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:O:79:TRP:CZ2	5:R:1:MET:N	2.66	0.56
7:T:17:THR:O	7:T:21:LEU:CB	2.52	0.56
2:B:155:LEU:HD23	2:B:155:LEU:C	2.25	0.56
2:B:155:LEU:HD23	2:B:155:LEU:O	2.04	0.56
1:A:88:TRP:HZ2	2:B:58:ASP:HB3	1.69	0.56
1:N:149:LYS:HD2	1:N:175:VAL:CG2	2.34	0.56
1:N:213:GLY:O	5:R:30:LYS:CD	2.49	0.56
4:Q:13:MET:SD	4:Q:13:MET:N	2.78	0.56
2:O:32:TRP:N	2:O:33:PRO:CD	2.67	0.56
4:Q:66:VAL:C	4:Q:68:LYS:H	2.09	0.56
1:A:26:VAL:HB	2:B:29:GLU:CG	2.35	0.56
2:B:41:PRO:O	2:B:45:MET:HG3	2.06	0.56
7:G:30:LYS:CB	7:G:31:ARG:NH2	2.66	0.56
5:E:9:ILE:HG23	6:F:12:SER:HB2	1.88	0.56
5:E:5:ALA:HB2	6:F:5:MET:HB2	1.87	0.56
2:O:141:ILE:O	2:O:144:GLY:N	2.39	0.56
3:C:185:LYS:HE2	3:C:195:TYR:CB	2.35	0.56
7:T:28:GLN:O	7:T:31:ARG:HB3	2.06	0.56
3:P:93:GLU:CA	3:P:97:GLU:HB2	2.28	0.56
1:A:106:LEU:HD12	5:E:18:ILE:HD12	1.85	0.56
3:P:219:VAL:HG11	3:P:223:GLN:NE2	2.19	0.56
1:N:109:GLY:C	1:N:111:LYS:H	2.08	0.56
1:N:118:TRP:CZ3	2:O:108:LEU:O	2.58	0.56
5:R:9:ILE:HG23	6:S:12:SER:HB2	1.88	0.56
5:E:6:VAL:O	5:E:10:VAL:HB	2.04	0.56
4:D:65:LYS:NZ	4:D:158:ASP:HB3	2.20	0.56
3:P:38:GLN:OE1	7:T:15:PHE:CE1	2.59	0.56
3:P:54:TYR:OH	3:P:58:LEU:HD22	2.05	0.56
3:P:12:THR:OG1	3:P:13:PRO:HD2	2.06	0.56
1:A:14:ILE:CD1	1:A:15:GLN:N	2.67	0.56
4:Q:73:HIS:O	4:Q:74:ASN:HB2	2.04	0.56
6:F:15:LEU:HD12	7:G:26:TYR:OH	2.05	0.56
7:T:11:LEU:HA	7:T:14:VAL:CG1	2.36	0.56
7:T:21:LEU:C	7:T:23:TYR:N	2.58	0.56
3:P:223:GLN:O	3:P:224:ALA:HB3	2.05	0.56
1:A:13:GLU:O	1:A:17:LEU:HD12	2.06	0.56
2:O:112:PRO:O	2:O:116:ASN:CB	2.53	0.56
3:C:271:MET:HG3	4:D:22:LEU:CD2	2.26	0.56
8:H:6:LEU:HA	8:H:9:VAL:HG12	1.87	0.56
2:O:86:GLN:HG2	2:O:143:LEU:HA	1.87	0.56
5:R:6:VAL:O	5:R:10:VAL:HB	2.05	0.56
2:O:32:TRP:N	2:O:33:PRO:HD3	2.20	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:37:LEU:HD23	2:B:40:PHE:CE2	2.41	0.56
7:G:25:ALA:O	7:G:29:TYR:CG	2.59	0.56
3:C:86:PRO:CG	3:C:89:ARG:HG3	2.36	0.56
1:N:15:GLN:O	1:N:17:LEU:N	2.39	0.56
1:A:77:SER:O	1:A:78:PHE:HB2	2.05	0.56
6:S:12:SER:O	6:S:16:ILE:HG12	2.05	0.56
4:Q:100:ARG:HD2	4:Q:102:TYR:CZ	2.40	0.56
7:G:11:LEU:CA	7:G:14:VAL:HG13	2.36	0.56
3:C:94:LEU:CD2	3:C:94:LEU:H	2.12	0.56
4:D:131:SER:CB	4:D:144:ALA:HA	2.36	0.56
4:Q:152:HIS:O	4:Q:153:ALA:HB2	2.05	0.56
2:B:36:LEU:O	2:B:39:VAL:HG22	2.06	0.56
5:R:7:PHE:CA	5:R:10:VAL:HG12	2.23	0.56
2:O:74:GLU:CG	2:O:75:ILE:H	2.18	0.56
2:O:96:LEU:O	2:O:100:LEU:HB2	2.05	0.56
1:A:72:ILE:O	1:A:79:GLY:HA3	2.06	0.56
3:C:178:GLY:HA2	3:C:225:VAL:HG13	1.86	0.56
4:Q:156:GLN:HB2	4:Q:161:VAL:HG11	1.88	0.56
12:B:307:OPC:CBI	12:B:307:OPC:HAP1	2.26	0.56
6:S:26:LEU:HD21	8:U:17:TRP:HE1	1.70	0.56
12:N:1306:OPC:HBQ2	12:N:1306:OPC:CAT	2.36	0.56
7:G:17:THR:O	7:G:21:LEU:CB	2.54	0.55
7:G:30:LYS:O	7:G:31:ARG:NE	2.39	0.55
3:P:134:PRO:CB	3:P:142:ILE:HD13	2.35	0.55
2:O:91:LEU:HD22	2:O:96:LEU:HD13	1.88	0.55
4:D:57:LYS:H	4:D:82:GLN:CG	2.20	0.55
4:Q:138:ARG:HH11	4:Q:138:ARG:HB3	1.69	0.55
3:C:269:GLN:HE22	7:G:24:ALA:CA	2.19	0.55
6:S:11:LEU:O	6:S:15:LEU:HD23	2.06	0.55
2:O:75:ILE:HD12	2:O:75:ILE:N	2.22	0.55
1:N:31:ASN:HB3	1:N:34:TYR:CD2	2.41	0.55
2:O:30:PRO:HG2	2:O:31:ALA:H	1.71	0.55
3:P:276:LYS:HA	3:P:276:LYS:HE3	1.89	0.55
5:E:26:ILE:HD12	5:E:26:ILE:N	2.22	0.55
12:Q:1307:OPC:HBZ2	12:Q:1307:OPC:CBF	2.37	0.55
3:P:255:VAL:O	3:P:258:MET:HG3	2.07	0.55
3:P:194:LYS:HD3	3:P:212:PRO:HA	1.89	0.55
8:U:3:ILE:HG13	8:U:4:ASP:OD1	2.06	0.55
1:A:166:SER:O	1:A:170:ARG:HG2	2.06	0.55
4:D:73:HIS:HE2	4:D:91:ILE:CG2	2.19	0.55
4:Q:155:VAL:HA	4:Q:159:ASN:O	2.05	0.55
3:C:218:ILE:CG1	3:C:232:THR:HG22	2.36	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:C:216:GLU:HB2	3:C:233:ASN:OD1	2.06	0.55
7:T:14:VAL:HG23	7:T:15:PHE:N	2.21	0.55
2:B:150:PRO:HG2	2:B:151:LEU:H	1.72	0.55
3:P:168:ASN:HA	3:P:172:PHE:CE1	2.40	0.55
3:P:26:HIS:ND1	3:P:154:ASN:OD1	2.39	0.55
4:Q:101:ASP:HA	4:Q:153:ALA:CB	2.37	0.55
2:B:31:ALA:O	2:B:32:TRP:CB	2.46	0.55
7:G:28:GLN:O	7:G:30:LYS:N	2.39	0.55
12:N:1306:OPC:HBQ2	12:N:1306:OPC:HAU2	1.89	0.55
1:A:108:GLY:HA2	1:A:110:PHE:CE2	2.41	0.55
1:N:28:PRO:O	2:O:30:PRO:HG2	2.07	0.55
2:O:32:TRP:CZ3	2:O:36:LEU:HG	2.42	0.55
5:R:27:LYS:O	5:R:31:LEU:HD22	2.06	0.55
4:D:168:THR:HA	4:D:176:PRO:CD	2.36	0.55
4:Q:92:VAL:HG21	4:Q:100:ARG:HB2	1.87	0.55
3:C:185:LYS:HE3	3:C:217:LEU:HD11	1.89	0.55
1:N:141:ASP:HA	2:O:65:PRO:CD	2.35	0.55
1:A:105:TYR:CD1	13:B:201:CLA:HMB1	2.41	0.55
1:A:88:TRP:CZ2	2:B:58:ASP:HB3	2.42	0.55
3:C:226:LYS:CG	3:C:229:GLU:HB3	2.36	0.55
3:C:111:ASP:O	3:C:114:LEU:HD21	2.06	0.55
1:N:213:GLY:O	5:R:26:ILE:HG23	2.06	0.55
4:Q:115:VAL:CG1	4:Q:124:PHE:HB3	2.35	0.55
1:A:95:LEU:HD11	5:E:10:VAL:HG11	1.88	0.55
2:B:149:LEU:N	2:B:149:LEU:CD1	2.66	0.55
3:P:119:LEU:H	3:P:119:LEU:HD12	1.69	0.55
1:N:193:TRP:O	1:N:197:VAL:HG23	2.07	0.55
3:P:82:PHE:O	3:P:112:ASN:HB3	2.07	0.55
5:R:5:ALA:HB2	6:S:5:MET:HB2	1.89	0.55
4:D:104:ILE:HG21	4:D:148:LEU:HB3	1.89	0.55
4:D:138:ARG:CZ	4:D:138:ARG:HB3	2.37	0.55
1:A:211:ILE:HD12	1:A:211:ILE:N	2.14	0.55
2:B:42:VAL:CG2	3:C:272:LEU:HD13	2.37	0.55
4:Q:104:ILE:HG21	4:Q:124:PHE:HE2	1.71	0.55
2:O:86:GLN:HG3	2:O:143:LEU:HD22	1.89	0.55
2:B:88:LEU:HD11	2:B:101:MET:SD	2.46	0.55
3:P:185:LYS:HD3	3:P:195:TYR:CD2	2.41	0.55
2:O:122:ASN:ND2	2:O:125:ARG:NH1	2.54	0.55
3:C:187:GLU:HG3	3:C:188:ASP:N	2.22	0.55
4:D:88:PRO:O	4:D:105:ASN:HA	2.07	0.55
3:C:40:VAL:HG13	3:C:247:ILE:HD11	1.89	0.55
3:P:275:LYS:NZ	3:P:276:LYS:NZ	2.55	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
12:D:306:OPC:CBM	12:D:306:OPC:HAR2	2.38	0.54
8:H:3:ILE:HG13	8:H:4:ASP:OD1	2.06	0.54
12:Q:1307:OPC:HBA1	12:Q:1307:OPC:HBP2	1.89	0.54
3:P:171:VAL:HG12	3:P:233:ASN:O	2.07	0.54
1:A:47:GLN:HG2	9:A:301:HEM:C3B	2.42	0.54
4:D:16:ARG:C	4:D:18:PHE:N	2.60	0.54
3:C:283:GLN:C	3:C:285:ALA:H	2.10	0.54
5:R:27:LYS:O	5:R:31:LEU:HB2	2.07	0.54
1:A:62:VAL:HG12	1:A:140:TRP:CH2	2.42	0.54
8:U:3:ILE:HG23	8:U:4:ASP:N	2.22	0.54
4:D:87:ASP:HB3	4:D:105:ASN:HB3	1.89	0.54
3:P:174:ALA:HB3	3:P:228:GLY:HA2	1.89	0.54
4:D:106:ALA:HB1	4:D:115:VAL:HB	1.89	0.54
4:Q:90:TYR:O	4:Q:91:ILE:HD13	2.08	0.54
12:Q:1307:OPC:HAP2	12:Q:1307:OPC:HAH2	1.88	0.54
2:O:106:LEU:CD2	13:O:1201:CLA:H151	2.36	0.54
1:N:141:ASP:C	2:O:64:GLU:HG2	2.27	0.54
3:P:199:ILE:HB	3:P:208:VAL:HA	1.90	0.54
3:P:171:VAL:HB	3:P:231:LEU:HD12	1.90	0.54
3:P:231:LEU:HD13	3:P:233:ASN:H	1.71	0.54
4:D:133:TYR:CD2	4:D:148:LEU:HG	2.42	0.54
4:Q:74:ASN:O	4:Q:94:GLU:O	2.25	0.54
4:Q:84:LEU:HD21	4:Q:164:PRO:HA	1.89	0.54
2:B:124:PHE:C	2:B:124:PHE:CD1	2.80	0.54
12:B:307:OPC:HAL2	12:B:307:OPC:CAP	2.37	0.54
3:C:273:ILE:O	3:C:278:GLN:HG2	2.08	0.54
2:O:79:TRP:HB3	5:R:7:PHE:HZ	1.73	0.54
1:A:33:PHE:HE1	5:E:14:LEU:HD22	1.71	0.54
3:C:20:ILE:HG21	3:C:75:VAL:HG11	1.88	0.54
3:P:275:LYS:O	3:P:279:VAL:HB	2.08	0.54
4:Q:73:HIS:NE2	4:Q:77:ASP:O	2.40	0.54
3:C:219:VAL:O	3:C:232:THR:HG21	2.07	0.54
3:P:231:LEU:HD22	3:P:232:THR:N	2.21	0.54
8:U:11:LEU:HB3	8:U:15:PHE:HZ	1.72	0.54
8:U:8:TRP:HA	8:U:11:LEU:HG	1.89	0.54
3:C:117:GLY:O	3:C:118:PRO:C	2.45	0.54
3:C:118:PRO:C	3:C:120:PRO:HD3	2.28	0.54
2:B:114:ILE:O	2:B:116:ASN:N	2.40	0.54
4:D:147:SER:CB	4:D:171:ARG:NH2	2.70	0.54
4:Q:121:GLU:O	4:Q:123:LYS:HG3	2.08	0.54
4:Q:81:VAL:O	4:Q:82:GLN:HG2	2.08	0.54
8:H:3:ILE:HG23	8:H:4:ASP:N	2.21	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:O:148:ALA:C	2:O:149:LEU:O	2.44	0.54
3:C:62:ALA:HB2	3:C:68:VAL:HG11	1.90	0.54
1:N:144:GLY:O	1:N:148:VAL:HG12	2.08	0.54
4:D:76:GLY:O	4:D:117:TRP:CH2	2.60	0.54
4:Q:157:ASP:O	4:Q:158:ASP:HB2	2.08	0.54
3:C:281:LYS:H	3:C:281:LYS:HD2	1.73	0.54
5:E:9:ILE:HG12	6:F:13:PHE:HB2	1.90	0.54
3:C:221:GLU:H	3:C:221:GLU:CD	2.12	0.54
3:P:79:PRO:HD3	3:P:149:ILE:HG12	1.89	0.54
1:A:14:ILE:CD1	1:A:15:GLN:H	2.18	0.54
4:Q:73:HIS:HA	4:Q:77:ASP:CG	2.28	0.54
7:G:30:LYS:HB3	7:G:31:ARG:HH22	1.71	0.54
7:T:22:PHE:O	7:T:26:TYR:CE1	2.61	0.54
3:C:226:LYS:HZ1	3:C:230:ALA:HB3	1.70	0.54
1:N:116:LEU:HD11	1:N:208:LYS:NZ	2.23	0.54
1:N:118:TRP:HD1	9:N:302:HEM:HMD2	1.72	0.54
4:Q:57:LYS:O	4:Q:64:VAL:HG23	2.08	0.54
5:E:3:LEU:O	5:E:4:GLY:C	2.47	0.54
3:C:194:LYS:HD3	3:C:212:PRO:HA	1.89	0.54
7:T:9:LEU:HG	7:T:10:VAL:H	1.72	0.54
3:P:178:GLY:O	3:P:225:VAL:HA	2.07	0.54
2:B:96:LEU:HD22	2:B:100:LEU:HG	1.90	0.54
3:P:283:GLN:C	3:P:285:ALA:N	2.62	0.54
2:O:39:VAL:O	2:O:42:VAL:HB	2.09	0.53
4:Q:132:GLN:NE2	4:Q:141:ARG:HB2	2.23	0.53
2:B:40:PHE:N	2:B:41:PRO:HD2	2.22	0.53
3:C:273:ILE:HG21	7:G:27:GLN:HB3	1.89	0.53
9:N:303:HEM:CBC	2:O:44:ILE:HD11	2.39	0.53
3:C:185:LYS:CG	3:C:195:TYR:HB3	2.30	0.53
2:B:146:GLY:O	2:B:147:ALA:HB3	2.08	0.53
8:U:10:ALA:O	8:U:14:VAL:HG23	2.08	0.53
1:A:166:SER:HB3	1:A:170:ARG:NH2	2.22	0.53
3:C:55:ASP:OD1	3:C:57:LYS:CG	2.56	0.53
4:Q:84:LEU:O	4:Q:85:LYS:HB2	2.07	0.53
8:H:5:VAL:O	8:H:9:VAL:HB	2.08	0.53
2:O:57:LEU:HD21	3:P:258:MET:HE1	1.89	0.53
6:S:15:LEU:HD12	7:T:26:TYR:OH	2.09	0.53
7:T:10:VAL:O	7:T:14:VAL:HG13	2.08	0.53
3:C:226:LYS:CE	3:C:229:GLU:HB3	2.37	0.53
3:C:54:TYR:OH	3:C:58:LEU:HD22	2.09	0.53
2:O:96:LEU:HD22	2:O:100:LEU:HG	1.89	0.53
2:B:48:PHE:CE2	2:B:52:VAL:HG21	2.43	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:R:18:ILE:HG22	5:R:22:ILE:HD11	1.91	0.53
4:Q:81:VAL:HB	4:Q:89:THR:CB	2.38	0.53
12:D:306:OPC:HBL1	12:D:306:OPC:CAR	2.28	0.53
3:P:216:GLU:O	3:P:233:ASN:CB	2.56	0.53
1:A:159:PRO:C	1:A:161:VAL:H	2.06	0.53
1:A:207:ARG:C	1:A:208:LYS:HD2	2.28	0.53
1:A:28:PRO:HD2	2:B:32:TRP:HB2	1.91	0.53
5:E:23:ILE:CG2	5:E:24:PHE:N	2.71	0.53
2:B:20:LYS:CG	2:B:21:GLY:N	2.55	0.53
5:R:1:MET:H1	5:R:1:MET:HE2	1.74	0.53
3:C:187:GLU:HG3	3:C:188:ASP:H	1.74	0.53
7:T:11:LEU:HA	7:T:14:VAL:HG13	1.91	0.53
10:A:304:TDS:HBC1	2:B:85:PHE:CD2	2.43	0.53
2:B:151:LEU:C	2:B:153:LYS:N	2.62	0.53
3:C:79:PRO:HD3	3:C:149:ILE:HG12	1.90	0.53
4:D:64:VAL:HG12	4:D:69:PHE:HE2	1.73	0.53
4:D:94:GLU:OE1	4:D:98:ALA:O	2.26	0.53
4:Q:109:THR:OG1	4:Q:145:PRO:HG2	2.09	0.53
4:Q:150:LEU:HD23	4:Q:178:TRP:CH2	2.44	0.53
4:Q:64:VAL:CG1	4:Q:91:ILE:HG12	2.36	0.53
4:D:19:MET:O	4:D:21:LEU:N	2.41	0.53
1:N:200:LEU:HD22	12:N:1306:OPC:HBC1	1.90	0.53
2:O:38:TYR:CG	3:P:276:LYS:HD3	2.43	0.53
5:R:23:ILE:CG2	5:R:24:PHE:N	2.71	0.53
4:D:73:HIS:HB3	4:D:93:VAL:HG12	1.90	0.53
2:B:72:PRO:HB3	4:Q:127:PRO:CG	2.39	0.53
4:Q:178:TRP:H	4:Q:178:TRP:HE3	1.57	0.53
3:C:275:LYS:O	3:C:279:VAL:HB	2.07	0.53
3:P:146:LYS:CD	3:P:246:GLU:HG3	2.29	0.53
1:N:142:GLN:HE21	2:O:64:GLU:CD	2.12	0.53
3:C:146:LYS:CD	3:C:246:GLU:HG3	2.32	0.53
3:C:119:LEU:HD12	3:C:119:LEU:H	1.72	0.53
2:O:21:GLY:C	2:O:24:HIS:CD2	2.82	0.53
3:C:40:VAL:HG11	3:C:46:PHE:CD2	2.44	0.53
4:Q:109:THR:CG2	4:Q:177:TRP:CH2	2.92	0.53
11:Q:1305:PL9:H301	12:Q:1307:OPC:HBU1	1.89	0.53
3:C:219:VAL:HG11	3:C:223:GLN:NE2	2.21	0.53
2:B:136:GLY:HA3	13:B:201:CLA:C2C	2.39	0.53
4:Q:15:ARG:HH12	4:Q:18:PHE:HE2	1.56	0.53
4:Q:136:THR:O	4:Q:171:ARG:NH2	2.41	0.53
3:C:93:GLU:O	3:C:94:LEU:C	2.48	0.53
1:N:83:ARG:NH1	9:N:301:HEM:O1A	2.41	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
12:D:306:OPC:CAY	12:D:306:OPC:HBC2	2.40	0.53
6:F:26:LEU:HD21	8:H:17:TRP:NE1	2.24	0.53
2:O:133:PHE:HA	13:O:1201:CLA:CMB	2.39	0.53
3:P:191:GLY:O	3:P:192:ASN:HB2	2.08	0.53
2:B:26:TYR:HB2	2:B:27:TYR:CE1	2.43	0.53
6:F:20:TRP:O	6:F:24:VAL:HG23	2.09	0.53
1:N:28:PRO:CG	2:O:32:TRP:HB3	2.39	0.52
1:A:211:ILE:CD1	1:A:211:ILE:N	2.70	0.52
2:O:79:TRP:HZ2	5:R:1:MET:N	2.03	0.52
3:C:194:LYS:HB3	3:C:196:GLN:HE21	1.74	0.52
3:P:185:LYS:HE3	3:P:217:LEU:HD11	1.90	0.52
4:Q:20:ASN:C	4:Q:22:LEU:N	2.60	0.52
4:D:134:ASP:CG	4:D:135:GLU:N	2.61	0.52
4:Q:142:GLY:HA2	4:Q:144:ALA:H	1.73	0.52
1:N:182:ARG:HB3	10:N:1304:TDS:HAA1	1.92	0.52
3:C:217:LEU:CA	3:C:232:THR:HB	2.38	0.52
1:N:61:THR:HG22	1:N:62:VAL:N	2.23	0.52
3:C:202:ASP:O	3:C:204:GLY:N	2.42	0.52
8:U:6:LEU:O	8:U:8:TRP:N	2.42	0.52
8:U:8:TRP:O	8:U:8:TRP:HE3	1.93	0.52
2:O:20:LYS:HG3	2:O:21:GLY:N	2.24	0.52
1:A:32:ILE:C	1:A:34:TYR:H	2.13	0.52
1:N:104:VAL:HA	1:N:107:THR:HG22	1.91	0.52
3:P:282:VAL:CG2	4:Q:15:ARG:HE	2.23	0.52
5:R:18:ILE:HB	5:R:22:ILE:HD12	1.92	0.52
4:Q:153:ALA:HA	4:Q:162:LEU:HD11	1.92	0.52
4:Q:92:VAL:O	4:Q:99:ILE:HA	2.09	0.52
12:B:307:OPC:HAL2	12:B:307:OPC:HAP1	1.91	0.52
2:B:36:LEU:C	2:B:38:TYR:H	2.11	0.52
7:G:14:VAL:HG23	7:G:15:PHE:N	2.25	0.52
1:N:142:GLN:HG2	2:O:64:GLU:CB	2.39	0.52
7:T:28:GLN:C	7:T:30:LYS:N	2.61	0.52
10:A:304:TDS:HAX2	13:B:201:CLA:H93	1.90	0.52
3:P:218:ILE:H	3:P:232:THR:CG2	2.23	0.52
4:D:138:ARG:NH2	4:D:147:SER:HB3	2.24	0.52
2:B:125:ARG:O	2:B:127:PRO:HD3	2.09	0.52
3:C:277:LYS:HB3	3:C:277:LYS:NZ	2.25	0.52
3:P:94:LEU:HD22	3:P:94:LEU:N	2.12	0.52
8:U:12:LEU:HA	8:U:15:PHE:CD1	2.44	0.52
3:P:226:LYS:HG2	3:P:229:GLU:CB	2.39	0.52
3:P:117:GLY:N	3:P:118:PRO:HD2	2.24	0.52
1:A:189:PHE:CE2	1:N:52:PHE:HB2	2.45	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:C:105:PRO:HB3	3:C:110:GLN:O	2.10	0.52
3:P:62:ALA:HB2	3:P:68:VAL:HG11	1.92	0.52
9:A:303:HEM:HBC1	11:A:305:PL9:H161	1.92	0.52
2:B:38:TYR:CD1	3:C:276:LYS:HD3	2.44	0.52
2:O:78:GLU:HG2	2:O:80:TYR:H	1.74	0.52
2:O:79:TRP:HB3	5:R:7:PHE:CZ	2.45	0.52
3:P:277:LYS:HB3	3:P:277:LYS:NZ	2.25	0.52
3:C:180:ILE:HG13	3:C:199:ILE:HA	1.91	0.52
1:N:23:SER:O	1:N:25:TYR:CE1	2.61	0.52
2:O:32:TRP:CE3	2:O:36:LEU:HG	2.44	0.52
4:Q:104:ILE:HD11	4:Q:136:THR:HA	1.90	0.52
4:Q:152:HIS:HB2	4:Q:162:LEU:HG	1.90	0.52
4:Q:154:THR:O	4:Q:160:ILE:HG23	2.08	0.52
4:D:19:MET:HG3	4:D:20:ASN:N	2.25	0.52
2:O:24:HIS:O	2:O:25:ASN:CB	2.57	0.52
2:O:75:ILE:HG22	2:O:77:PRO:HD3	1.91	0.52
5:R:26:ILE:N	5:R:26:ILE:HD12	2.24	0.52
4:Q:100:ARG:HD2	4:Q:102:TYR:HH	1.74	0.52
4:Q:76:GLY:CA	4:Q:93:VAL:HG13	2.40	0.52
8:H:8:TRP:O	8:H:8:TRP:HE3	1.92	0.52
3:C:231:LEU:HD12	3:C:231:LEU:H	1.75	0.52
9:A:301:HEM:HBD1	9:A:301:HEM:HHA	1.92	0.52
3:P:124:TYR:HB3	3:P:127:ILE:HG12	1.92	0.52
1:N:209:GLN:HB3	2:O:27:TYR:HB2	1.92	0.52
4:D:73:HIS:CE1	4:D:91:ILE:O	2.63	0.52
4:Q:78:ARG:NH1	4:Q:100:ARG:HH21	2.06	0.52
4:Q:64:VAL:HG13	4:Q:69:PHE:CD2	2.45	0.52
8:H:6:LEU:O	8:H:8:TRP:N	2.42	0.52
11:Q:1305:PL9:H302	12:Q:1307:OPC:CBB	2.39	0.52
11:Q:1305:PL9:H201	12:Q:1307:OPC:OCC	2.08	0.52
3:C:28:ALA:HB3	3:C:238:GLY:O	2.10	0.52
3:P:218:ILE:HG13	3:P:232:THR:HG22	1.90	0.52
3:P:233:ASN:ND2	3:P:234:ASN:H	2.08	0.52
2:O:127:PRO:C	2:O:129:ALA:N	2.63	0.52
8:U:6:LEU:C	8:U:8:TRP:H	2.13	0.52
1:A:41:LEU:O	1:A:44:PHE:HB3	2.09	0.52
8:U:22:VAL:O	8:U:26:ARG:HG3	2.10	0.52
3:C:11:PRO:HA	3:C:107:LYS:NZ	2.25	0.52
4:D:116:PRO:CD	4:D:126:CYS:HA	2.40	0.52
4:D:55:THR:OG1	4:D:57:LYS:NZ	2.36	0.52
2:B:72:PRO:HB3	4:Q:127:PRO:HG2	1.90	0.52
5:E:21:GLY:O	5:E:25:ALA:HB2	2.10	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
12:Q:1307:OPC:HBX2	12:Q:1307:OPC:CBE	2.39	0.52
2:O:85:PHE:O	2:O:89:ARG:HG2	2.10	0.52
7:T:25:ALA:CB	7:T:29:TYR:OH	2.57	0.52
3:C:91:PRO:HD2	3:C:95:LYS:HG3	1.92	0.52
1:A:44:PHE:HD1	9:A:301:HEM:CBB	2.21	0.52
3:C:201:THR:O	3:C:203:SER:N	2.43	0.52
2:O:122:ASN:OD1	5:R:26:ILE:HG21	2.10	0.52
6:S:17:PHE:N	6:S:17:PHE:CD1	2.78	0.52
4:Q:84:LEU:HD13	4:Q:164:PRO:HD3	1.91	0.52
2:B:79:TRP:CG	2:B:80:TYR:N	2.78	0.52
3:P:185:LYS:CD	3:P:195:TYR:HD2	2.23	0.52
2:O:95:LEU:N	2:O:95:LEU:CD2	2.67	0.52
3:P:52:ILE:HG21	3:P:154:ASN:O	2.10	0.52
4:D:84:LEU:H	4:D:89:THR:HG23	1.75	0.52
2:O:24:HIS:O	2:O:25:ASN:HB3	2.10	0.52
4:Q:116:PRO:HD3	4:Q:127:PRO:HD3	1.92	0.51
7:G:15:PHE:C	7:G:17:THR:N	2.63	0.51
1:A:33:PHE:CZ	5:E:17:GLY:HA3	2.45	0.51
2:B:150:PRO:O	2:B:153:LYS:NZ	2.43	0.51
13:B:201:CLA:HHC	13:B:201:CLA:HBB1	1.93	0.51
5:E:18:ILE:HG22	5:E:22:ILE:HD11	1.92	0.51
3:P:202:ASP:O	3:P:204:GLY:N	2.44	0.51
3:P:55:ASP:OD1	3:P:57:LYS:HG3	2.10	0.51
3:C:59:GLN:HB2	3:C:67:LYS:HB3	1.93	0.51
1:N:114:ARG:NH2	1:N:209:GLN:H	2.08	0.51
1:N:29:HIS:O	1:N:30:VAL:CG1	2.45	0.51
1:N:106:LEU:HD12	5:R:18:ILE:CD1	2.40	0.51
4:D:78:ARG:HG3	4:D:117:TRP:CG	2.45	0.51
3:C:218:ILE:H	3:C:232:THR:CG2	2.22	0.51
3:P:221:GLU:CD	3:P:221:GLU:H	2.14	0.51
1:N:110:PHE:CD1	2:O:112:PRO:HB3	2.42	0.51
4:Q:150:LEU:HD11	4:Q:171:ARG:NH2	2.22	0.51
1:A:172:GLY:O	1:A:173:SER:C	2.48	0.51
2:B:39:VAL:O	2:B:42:VAL:HB	2.10	0.51
1:A:39:ILE:HD11	2:B:43:VAL:CG1	2.40	0.51
15:E:101:BCR:H362	6:F:17:PHE:HZ	1.75	0.51
5:E:20:VAL:CG1	5:E:21:GLY:N	2.69	0.51
2:O:143:LEU:O	2:O:145:ILE:N	2.43	0.51
2:B:79:TRP:HA	2:B:82:TYR:CD2	2.46	0.51
3:C:102:TYR:N	3:C:118:PRO:HG3	2.22	0.51
3:C:134:PRO:CB	3:C:142:ILE:HD13	2.39	0.51
3:C:50:VAL:O	3:C:50:VAL:HG12	2.10	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:P:50:VAL:HG13	3:P:151:LEU:HD21	1.92	0.51
1:A:75:GLU:HA	1:A:75:GLU:OE1	2.11	0.51
4:D:115:VAL:HG21	4:D:148:LEU:CD2	2.41	0.51
8:U:17:TRP:O	8:U:20:ALA:HB3	2.10	0.51
2:B:150:PRO:O	2:B:153:LYS:HD2	2.09	0.51
3:P:118:PRO:O	3:P:120:PRO:HD3	2.11	0.51
2:B:96:LEU:O	2:B:100:LEU:HB2	2.11	0.51
1:A:99:LEU:HD21	5:E:11:PHE:HB2	1.92	0.51
2:O:118:ASN:ND2	2:O:121:GLN:NE2	2.37	0.51
4:Q:88:PRO:CD	4:Q:107:VAL:HG23	2.41	0.51
4:Q:137:GLY:C	4:Q:171:ARG:HH12	2.14	0.51
4:Q:77:ASP:N	4:Q:93:VAL:HG12	2.18	0.51
4:D:155:VAL:HG12	4:D:156:GLN:N	2.25	0.51
1:N:26:VAL:HG23	1:N:26:VAL:O	2.11	0.51
5:R:23:ILE:O	5:R:27:LYS:CB	2.58	0.51
4:Q:133:TYR:CD2	4:Q:139:VAL:HB	2.46	0.51
4:Q:152:HIS:CA	4:Q:162:LEU:HG	2.41	0.51
2:B:32:TRP:H	2:B:33:PRO:CD	2.24	0.51
5:E:23:ILE:HG23	5:E:24:PHE:HD2	1.75	0.51
4:Q:35:LEU:O	4:Q:39:VAL:HG23	2.11	0.51
3:C:178:GLY:O	3:C:225:VAL:HA	2.11	0.51
4:Q:109:THR:HG22	4:Q:177:TRP:CH2	2.45	0.51
5:R:3:LEU:O	5:R:4:GLY:C	2.49	0.51
1:A:103:ARG:NH1	1:A:103:ARG:HG3	2.25	0.51
3:P:3:PHE:CD1	3:P:3:PHE:N	2.78	0.51
1:A:81:LEU:C	1:A:81:LEU:HD13	2.30	0.51
4:D:35:LEU:HD12	4:D:35:LEU:O	2.11	0.51
3:P:201:THR:O	3:P:203:SER:N	2.43	0.51
1:N:29:HIS:HE1	2:O:29:GLU:OE2	1.94	0.51
4:Q:16:ARG:C	4:Q:18:PHE:N	2.64	0.51
4:D:90:TYR:HD2	4:D:104:ILE:O	1.93	0.51
4:D:100:ARG:C	4:D:153:ALA:HB2	2.31	0.51
4:Q:152:HIS:HB2	4:Q:162:LEU:CB	2.41	0.51
4:Q:78:ARG:N	4:Q:78:ARG:HD3	2.25	0.51
3:C:262:ILE:HG12	7:G:17:THR:HG22	1.92	0.51
1:A:118:TRP:CZ3	2:B:108:LEU:O	2.64	0.51
2:O:74:GLU:CG	2:O:75:ILE:N	2.73	0.51
3:P:20:ILE:HG21	3:P:75:VAL:HG11	1.91	0.51
4:Q:152:HIS:CB	4:Q:162:LEU:HG	2.40	0.51
4:Q:162:LEU:O	4:Q:163:THR:HB	2.11	0.51
5:E:16:PHE:O	5:E:20:VAL:HB	2.11	0.51
3:C:266:MET:HE3	7:G:23:TYR:CB	2.41	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:98:VAL:HA	2:B:101:MET:HE2	1.93	0.51
8:U:5:VAL:O	8:U:9:VAL:HB	2.11	0.51
3:P:226:LYS:CD	3:P:229:GLU:HB3	2.40	0.51
1:N:14:ILE:HG13	1:N:14:ILE:O	2.11	0.51
4:D:115:VAL:HG12	4:D:124:PHE:HB3	1.92	0.50
1:N:95:LEU:HG	5:R:7:PHE:HB3	1.94	0.50
2:O:78:GLU:OE1	2:O:80:TYR:HB2	2.11	0.50
2:B:79:TRP:C	2:B:81:LEU:H	2.14	0.50
3:P:187:GLU:HG3	3:P:188:ASP:N	2.26	0.50
3:C:61:VAL:O	3:C:157:ARG:NH1	2.42	0.50
1:N:150:ILE:HD12	2:O:75:ILE:HG12	1.93	0.50
2:O:91:LEU:O	2:O:91:LEU:HD12	2.11	0.50
4:D:95:SER:C	4:D:97:GLU:H	2.14	0.50
5:R:5:ALA:HB1	6:S:9:ALA:CB	2.21	0.50
4:Q:123:LYS:O	4:Q:125:LYS:N	2.44	0.50
4:Q:132:GLN:NE2	4:Q:141:ARG:HD3	2.22	0.50
4:Q:149:ALA:O	4:Q:150:LEU:HD13	2.11	0.50
2:B:127:PRO:C	2:B:129:ALA:N	2.63	0.50
2:O:57:LEU:O	3:P:38:GLN:NE2	2.45	0.50
8:U:24:TRP:CD2	8:U:25:GLY:N	2.79	0.50
3:P:194:LYS:HB3	3:P:196:GLN:HE21	1.75	0.50
3:P:231:LEU:CD1	3:P:231:LEU:H	2.24	0.50
4:D:59:LYS:HB3	4:D:80:LEU:O	2.10	0.50
3:C:53:PRO:O	3:C:54:TYR:HB2	2.10	0.50
3:C:34:VAL:HG22	3:C:151:LEU:HD22	1.93	0.50
2:O:73:LEU:C	2:O:73:LEU:HD12	2.31	0.50
1:N:214:PRO:HA	5:R:29:ILE:O	2.12	0.50
4:D:118:ASN:OD1	4:D:123:LYS:C	2.50	0.50
4:D:78:ARG:NE	4:D:92:VAL:HG11	2.23	0.50
4:Q:70:LEU:HD22	4:Q:71:GLU:OE1	2.12	0.50
7:G:26:TYR:HD1	7:G:26:TYR:H	1.59	0.50
1:N:140:TRP:O	2:O:65:PRO:HG3	2.11	0.50
2:B:142:TRP:CD1	2:B:155:LEU:C	2.85	0.50
2:B:142:TRP:NE1	2:B:154:THR:O	2.41	0.50
3:C:3:PHE:CD1	3:C:3:PHE:N	2.78	0.50
3:C:134:PRO:HB3	3:C:142:ILE:CD1	2.40	0.50
1:N:31:ASN:O	1:N:32:ILE:HB	2.12	0.50
4:Q:156:GLN:O	4:Q:157:ASP:CB	2.59	0.50
4:Q:94:GLU:HB3	4:Q:100:ARG:CD	2.12	0.50
7:G:21:LEU:C	7:G:23:TYR:H	2.13	0.50
1:N:38:GLY:HA3	9:N:303:HEM:C4C	2.46	0.50
2:O:76:LEU:HD23	2:O:76:LEU:H	1.75	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:C:231:LEU:CD1	3:C:232:THR:H	2.23	0.50
7:T:22:PHE:C	7:T:26:TYR:CE1	2.85	0.50
8:U:9:VAL:HG22	8:U:13:VAL:CG2	2.41	0.50
3:C:199:ILE:HG22	3:C:200:GLN:N	2.26	0.50
1:N:33:PHE:O	1:N:36:LEU:HD23	2.11	0.50
4:D:106:ALA:HB3	4:D:114:VAL:HG13	1.94	0.50
4:Q:147:SER:HB3	4:Q:171:ARG:NH1	2.26	0.50
15:E:101:BCR:H362	6:F:17:PHE:CZ	2.46	0.50
3:C:231:LEU:HD22	3:C:232:THR:H	1.77	0.50
3:C:233:ASN:ND2	3:C:234:ASN:H	2.09	0.50
2:O:36:LEU:O	2:O:39:VAL:HG22	2.11	0.50
4:Q:116:PRO:CD	4:Q:127:PRO:HD3	2.41	0.50
4:Q:156:GLN:HB2	4:Q:161:VAL:CG1	2.42	0.50
4:D:68:LYS:H	4:D:68:LYS:HD2	1.75	0.50
2:B:79:TRP:C	2:B:81:LEU:N	2.64	0.50
2:O:22:MET:O	2:O:24:HIS:N	2.44	0.50
2:B:86:GLN:HG3	2:B:143:LEU:HD22	1.94	0.50
1:A:14:ILE:HA	1:A:17:LEU:HD12	1.94	0.50
4:D:110:HIS:HE1	4:D:129:HIS:HB2	1.77	0.50
4:Q:123:LYS:HA	4:Q:133:TYR:O	2.11	0.50
12:D:306:OPC:CCA	12:D:306:OPC:HBW2	2.39	0.50
1:A:158:ILE:HB	1:A:162:GLY:HA2	1.91	0.50
3:C:178:GLY:C	3:C:225:VAL:HA	2.32	0.50
4:D:167:GLU:O	4:D:176:PRO:HB3	2.11	0.50
4:D:167:GLU:O	4:D:176:PRO:CB	2.59	0.50
4:D:149:ALA:HA	4:D:178:TRP:CE2	2.47	0.50
4:Q:134:ASP:C	4:Q:136:THR:H	2.14	0.50
4:Q:80:LEU:HD23	4:Q:80:LEU:C	2.32	0.50
1:A:27:PRO:HG3	2:B:33:PRO:CG	2.41	0.50
3:C:274:LEU:HD11	8:H:23:VAL:HG21	1.92	0.50
1:N:95:LEU:HD13	1:N:96:MET:HE3	1.92	0.50
3:P:86:PRO:CG	3:P:89:ARG:HG3	2.41	0.50
2:B:79:TRP:CD1	2:B:80:TYR:N	2.80	0.50
3:P:173:THR:HB	3:P:229:GLU:HA	1.92	0.50
3:C:52:ILE:HG12	3:C:153:ALA:CB	2.40	0.50
3:C:199:ILE:HB	3:C:208:VAL:HA	1.94	0.50
3:C:56:THR:HG22	3:C:122:GLU:HB3	1.92	0.50
4:D:121:GLU:CD	4:D:123:LYS:HE2	2.32	0.50
4:Q:100:ARG:CD	4:Q:102:TYR:OH	2.57	0.50
4:Q:56:ALA:HA	4:Q:81:VAL:CG1	2.25	0.50
9:N:303:HEM:O2D	11:Q:1305:PL9:H532	2.11	0.50
5:R:1:MET:H1	5:R:1:MET:CE	2.24	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:N:41:LEU:O	1:N:44:PHE:HB3	2.12	0.50
1:N:115:GLU:O	1:N:119:ILE:HG13	2.12	0.49
4:D:99:ILE:HD12	4:D:153:ALA:HB3	1.94	0.49
4:Q:64:VAL:HA	4:Q:69:PHE:CE1	2.47	0.49
3:C:269:GLN:HE22	7:G:24:ALA:HA	1.75	0.49
2:B:32:TRP:CH2	3:C:277:LYS:HD2	2.43	0.49
3:C:277:LYS:HE3	7:G:31:ARG:CG	2.42	0.49
6:F:17:PHE:N	6:F:17:PHE:CD1	2.80	0.49
1:A:177:GLN:H	1:A:177:GLN:CD	2.13	0.49
3:P:3:PHE:CZ	3:P:119:LEU:HD11	2.41	0.49
3:P:134:PRO:HB3	3:P:142:ILE:CD1	2.36	0.49
4:Q:120:ALA:C	4:Q:122:ASN:H	2.16	0.49
2:B:117:VAL:HG22	2:B:117:VAL:O	2.11	0.49
5:R:9:ILE:HB	6:S:9:ALA:HB1	1.94	0.49
4:D:165:TRP:HA	4:D:167:GLU:OE2	2.12	0.49
4:Q:99:ILE:HD12	4:Q:160:ILE:HD13	1.94	0.49
2:B:127:PRO:O	2:B:129:ALA:N	2.42	0.49
11:Q:1305:PL9:C30	12:Q:1307:OPC:HBU1	2.42	0.49
1:A:36:LEU:CD2	1:A:103:ARG:HB2	2.43	0.49
2:B:58:ASP:OD1	3:C:146:LYS:NZ	2.45	0.49
4:D:59:LYS:CD	4:D:79:VAL:HB	2.42	0.49
2:O:22:MET:C	2:O:24:HIS:H	2.14	0.49
5:E:30:LYS:CB	5:E:30:LYS:NZ	2.75	0.49
1:N:106:LEU:HD12	5:R:18:ILE:HD12	1.95	0.49
4:Q:141:ARG:HG2	4:Q:141:ARG:HH11	1.77	0.49
4:Q:154:THR:O	4:Q:160:ILE:CA	2.58	0.49
7:G:11:LEU:C	7:G:14:VAL:HG13	2.33	0.49
3:P:91:PRO:HD2	3:P:95:LYS:HG3	1.94	0.49
3:P:93:GLU:O	3:P:94:LEU:C	2.49	0.49
1:A:91:SER:CB	2:B:79:TRP:HZ3	2.25	0.49
3:C:226:LYS:O	3:C:227:ALA:HB2	2.12	0.49
3:C:26:HIS:ND1	3:C:154:ASN:OD1	2.46	0.49
4:Q:170:PHE:HD1	4:Q:170:PHE:H	1.60	0.49
1:N:212:SER:HB2	2:O:122:ASN:OD1	2.12	0.49
4:D:102:TYR:CE2	4:D:136:THR:HG23	2.47	0.49
4:D:104:ILE:HG23	4:D:149:ALA:O	2.13	0.49
7:G:15:PHE:C	7:G:17:THR:H	2.15	0.49
2:B:22:MET:SD	2:B:22:MET:O	2.70	0.49
3:C:124:TYR:HB3	3:C:127:ILE:HG12	1.95	0.49
3:P:59:GLN:HB2	3:P:67:LYS:HB3	1.94	0.49
3:C:262:ILE:HG13	7:G:21:LEU:CG	2.27	0.49
6:S:7:TYR:O	6:S:11:LEU:HD12	2.12	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:147:ALA:HB1	2:B:152:ASP:HA	1.95	0.49
8:U:6:LEU:O	8:U:10:ALA:HB3	2.12	0.49
1:A:66:TYR:CZ	2:B:65:PRO:HD3	2.47	0.49
1:N:71:TYR:O	1:N:76:VAL:HG23	2.12	0.49
4:D:126:CYS:N	4:D:131:SER:O	2.38	0.49
4:Q:94:GLU:OE2	4:Q:100:ARG:HG2	2.12	0.49
1:A:114:ARG:NH1	1:A:208:LYS:HB3	2.28	0.49
2:B:32:TRP:N	2:B:33:PRO:CD	2.76	0.49
2:O:79:TRP:CH2	5:R:1:MET:N	2.79	0.49
2:O:64:GLU:CG	2:O:65:PRO:N	2.75	0.49
12:N:1306:OPC:HBQ2	12:N:1306:OPC:CAU	2.42	0.49
3:P:281:LYS:HA	3:P:284:ALA:CB	2.25	0.49
2:B:82:TYR:N	2:B:83:PRO:HD2	2.28	0.49
3:P:199:ILE:HG22	3:P:200:GLN:N	2.27	0.49
3:P:171:VAL:HG13	3:P:234:ASN:HB2	1.93	0.49
4:D:84:LEU:O	4:D:87:ASP:N	2.46	0.49
3:P:270:LEU:HG	3:P:274:LEU:CD1	2.43	0.49
3:C:174:ALA:HB3	3:C:228:GLY:HA2	1.93	0.49
1:N:111:LYS:HB2	1:N:114:ARG:NH1	2.26	0.49
4:Q:169:ASP:HB3	4:Q:172:THR:OG1	2.13	0.49
5:E:26:ILE:HG13	5:E:29:ILE:HD12	1.95	0.49
7:G:21:LEU:C	7:G:23:TYR:N	2.63	0.49
3:C:191:GLY:O	3:C:192:ASN:HB2	2.13	0.49
3:C:217:LEU:HA	3:C:232:THR:CB	2.39	0.49
1:A:97:MET:O	1:A:101:VAL:HG23	2.13	0.49
2:B:141:ILE:O	2:B:144:GLY:N	2.45	0.49
3:P:275:LYS:HB2	4:Q:19:MET:CE	2.43	0.49
3:P:275:LYS:NZ	3:P:276:LYS:HZ2	2.10	0.49
2:B:118:ASN:O	2:B:121:GLN:O	2.30	0.49
2:O:147:ALA:O	2:O:149:LEU:N	2.45	0.49
3:P:226:LYS:O	3:P:227:ALA:HB2	2.12	0.49
1:A:142:GLN:HE22	2:B:67:ASN:H	1.59	0.49
4:D:174:GLU:HG2	4:D:175:LYS:O	2.13	0.49
4:Q:163:THR:O	4:Q:164:PRO:O	2.31	0.49
2:B:32:TRP:N	2:B:33:PRO:HD3	2.28	0.49
4:D:19:MET:C	4:D:21:LEU:H	2.14	0.49
2:O:146:GLY:O	2:O:149:LEU:CB	2.61	0.49
3:P:26:HIS:ND1	3:P:154:ASN:ND2	2.61	0.49
1:A:178:ALA:O	1:A:182:ARG:HG3	2.13	0.49
2:B:102:ALA:O	2:B:105:PRO:HD2	2.12	0.49
4:D:166:THR:CA	4:D:179:VAL:HG22	2.43	0.49
1:A:114:ARG:HB3	1:A:208:LYS:NZ	2.27	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:D:19:MET:C	4:D:21:LEU:N	2.65	0.49
7:G:25:ALA:CB	7:G:29:TYR:OH	2.61	0.49
1:A:102:PHE:O	1:A:103:ARG:C	2.50	0.49
2:B:79:TRP:CD1	2:B:79:TRP:C	2.86	0.49
3:C:182:LYS:CG	3:C:198:SER:HB2	2.40	0.49
3:C:199:ILE:HB	3:C:207:VAL:O	2.13	0.49
4:D:35:LEU:O	4:D:39:VAL:HG23	2.13	0.49
2:O:153:LYS:O	2:O:155:LEU:N	2.40	0.49
1:N:110:PHE:O	1:N:111:LYS:O	2.31	0.48
2:O:18:LEU:HD13	2:O:31:ALA:HB2	1.93	0.48
4:D:54:THR:HG23	4:D:82:GLN:HG3	1.95	0.48
5:E:26:ILE:HG22	5:E:26:ILE:O	2.12	0.48
1:N:97:MET:O	1:N:101:VAL:HG23	2.12	0.48
2:O:59:PRO:HD2	3:P:248:VAL:HG21	1.95	0.48
3:P:9:TYR:C	3:P:11:PRO:HD2	2.33	0.48
3:C:9:TYR:C	3:C:11:PRO:HD2	2.33	0.48
1:A:99:LEU:CD2	5:E:11:PHE:HB2	2.43	0.48
3:C:98:VAL:O	3:C:99:GLY:O	2.31	0.48
4:Q:102:TYR:H	4:Q:153:ALA:HB2	1.78	0.48
7:G:17:THR:O	7:G:21:LEU:HB2	2.13	0.48
7:G:21:LEU:O	7:G:21:LEU:HD13	2.12	0.48
7:G:27:GLN:HA	7:G:27:GLN:NE2	2.28	0.48
2:O:79:TRP:HA	2:O:82:TYR:CZ	2.48	0.48
6:S:26:LEU:HD22	8:U:17:TRP:HE1	1.75	0.48
3:P:219:VAL:O	3:P:232:THR:HG21	2.13	0.48
8:U:8:TRP:O	8:U:11:LEU:HB2	2.13	0.48
3:P:226:LYS:CG	3:P:229:GLU:HB3	2.43	0.48
2:O:46:GLY:HA3	7:T:24:ALA:HB1	1.95	0.48
1:N:135:GLY:HA2	1:N:138:LEU:HG	1.96	0.48
4:Q:28:THR:OG1	11:Q:1305:PL9:H23	2.13	0.48
5:R:9:ILE:CG1	6:S:13:PHE:HB2	2.43	0.48
4:D:139:VAL:HG23	4:D:144:ALA:HB3	1.96	0.48
4:Q:88:PRO:O	4:Q:89:THR:HG23	2.13	0.48
4:Q:74:ASN:N	4:Q:93:VAL:HG11	2.13	0.48
8:H:7:GLY:O	8:H:11:LEU:HD11	2.13	0.48
4:D:66:VAL:CG2	4:D:158:ASP:C	2.82	0.48
1:A:104:VAL:O	1:A:107:THR:HG22	2.14	0.48
2:B:91:LEU:O	2:B:91:LEU:HD12	2.13	0.48
2:B:135:PHE:O	2:B:138:LEU:HB2	2.13	0.48
3:P:34:VAL:HG22	3:P:151:LEU:HD22	1.95	0.48
3:C:12:THR:OG1	3:C:13:PRO:HD2	2.12	0.48
1:N:104:VAL:O	1:N:107:THR:HG22	2.12	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:D:147:SER:HB2	4:D:171:ARG:NH2	2.28	0.48
3:C:266:MET:HG3	7:G:21:LEU:HD23	1.95	0.48
2:O:82:TYR:N	2:O:83:PRO:HD2	2.28	0.48
8:U:21:MET:O	8:U:24:TRP:HE3	1.96	0.48
4:D:107:VAL:HG12	4:D:109:THR:HG23	1.94	0.48
4:D:130:GLY:C	4:D:143:PRO:HG2	2.34	0.48
1:N:141:ASP:O	1:N:145:TYR:HB3	2.13	0.48
1:N:66:TYR:CG	2:O:65:PRO:HG2	2.48	0.48
3:P:273:ILE:O	3:P:278:GLN:HG2	2.13	0.48
3:C:91:PRO:HD2	3:C:95:LYS:HG2	1.94	0.48
3:P:231:LEU:CD1	3:P:231:LEU:N	2.77	0.48
1:N:109:GLY:HA3	9:N:302:HEM:HBD2	1.96	0.48
1:N:118:TRP:CH2	2:O:108:LEU:HB3	2.49	0.48
4:D:116:PRO:HD2	4:D:126:CYS:SG	2.54	0.48
4:Q:159:ASN:O	4:Q:161:VAL:N	2.46	0.48
4:Q:66:VAL:O	4:Q:70:LEU:N	2.46	0.48
1:A:35:CYS:HA	9:A:303:HEM:C3B	2.49	0.48
1:N:72:ILE:O	1:N:79:GLY:HA3	2.13	0.48
3:P:104:GLN:HG2	3:P:115:LEU:HB2	1.96	0.48
1:N:107:THR:O	2:O:123:PRO:HG2	2.13	0.48
2:O:18:LEU:CB	2:O:31:ALA:HB2	2.43	0.48
5:R:25:ALA:C	5:R:27:LYS:H	2.17	0.48
4:D:138:ARG:HH11	4:D:138:ARG:HB3	1.75	0.48
4:Q:115:VAL:HG13	4:Q:126:CYS:CA	2.43	0.48
4:Q:155:VAL:HG12	4:Q:156:GLN:H	1.79	0.48
4:Q:77:ASP:OD1	4:Q:93:VAL:HG11	2.14	0.48
3:C:41:LEU:HD21	7:G:11:LEU:HD21	1.96	0.48
4:D:15:ARG:HH22	4:D:17:GLN:HG2	1.75	0.48
8:H:24:TRP:CD2	8:H:25:GLY:N	2.82	0.48
2:O:58:ASP:OD1	5:R:2:ILE:HG21	2.13	0.48
1:N:99:LEU:HD21	5:R:11:PHE:HB2	1.96	0.48
3:C:1:TYR:CD2	3:C:119:LEU:HD21	2.46	0.48
3:C:251:ASP:HB3	3:C:254:ARG:HD2	1.94	0.48
1:N:104:VAL:HG11	9:N:302:HEM:HMD2	1.96	0.48
4:D:138:ARG:HE	4:D:171:ARG:HD2	1.77	0.48
4:Q:92:VAL:HG23	4:Q:100:ARG:CB	2.44	0.48
4:Q:94:GLU:CG	4:Q:95:SER:N	2.76	0.48
3:C:275:LYS:HE2	4:D:17:GLN:HB2	1.95	0.48
7:G:14:VAL:HG22	7:G:15:PHE:H	1.78	0.48
5:E:5:ALA:HB2	6:F:5:MET:CB	2.43	0.48
6:S:22:LEU:O	6:S:26:LEU:HD23	2.14	0.48
3:P:226:LYS:CE	3:P:229:GLU:HB3	2.44	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:N:58:TYR:HE2	1:N:60:PRO:HB3	1.79	0.48
4:D:61:GLY:HA2	2:O:66:ALA:O	2.14	0.48
2:O:36:LEU:C	2:O:38:TYR:H	2.17	0.48
4:D:118:ASN:HB2	4:D:122:ASN:H	1.78	0.48
4:D:138:ARG:HH21	4:D:147:SER:HB3	1.79	0.48
1:A:23:SER:O	1:A:25:TYR:HD1	1.97	0.48
1:A:30:VAL:O	1:A:30:VAL:HG23	2.14	0.48
1:N:191:LEU:HD11	10:N:1304:TDS:OBD	2.14	0.48
2:B:85:PHE:O	2:B:89:ARG:HG2	2.14	0.48
3:P:187:GLU:HG3	3:P:188:ASP:H	1.79	0.48
3:P:117:GLY:O	3:P:118:PRO:C	2.51	0.48
2:O:153:LYS:C	2:O:155:LEU:H	2.17	0.48
1:N:112:LYS:CB	1:N:113:PRO:HD3	2.31	0.47
3:P:275:LYS:HG3	4:Q:20:ASN:CA	2.43	0.47
5:R:26:ILE:O	5:R:26:ILE:HG22	2.12	0.47
4:D:94:GLU:N	4:D:94:GLU:OE1	2.40	0.47
4:D:68:LYS:N	4:D:68:LYS:HD2	2.29	0.47
4:D:70:LEU:HD23	4:D:70:LEU:O	2.14	0.47
7:T:14:VAL:HG22	7:T:15:PHE:H	1.79	0.47
3:P:218:ILE:CG1	3:P:232:THR:HG22	2.43	0.47
1:N:163:VAL:CG1	1:N:164:LEU:N	2.77	0.47
2:O:18:LEU:HD21	7:T:35:LEU:O	2.14	0.47
4:D:147:SER:HB3	4:D:171:ARG:NH2	2.29	0.47
4:Q:93:VAL:HG22	4:Q:94:GLU:H	1.79	0.47
2:B:126:ARG:O	2:B:129:ALA:HB3	2.14	0.47
2:O:79:TRP:CH2	5:R:4:GLY:CA	2.92	0.47
4:D:66:VAL:CG2	4:D:160:ILE:N	2.77	0.47
1:A:158:ILE:CB	1:A:162:GLY:HA3	2.40	0.47
1:N:20:ASP:O	1:N:21:VAL:O	2.32	0.47
3:P:47:LYS:HG3	3:P:128:VAL:HG13	1.96	0.47
1:N:34:TYR:CE2	1:N:103:ARG:HD3	2.48	0.47
3:P:279:VAL:O	3:P:279:VAL:HG12	2.14	0.47
3:P:275:LYS:HB2	4:Q:19:MET:HE2	1.95	0.47
2:B:38:TYR:CG	3:C:276:LYS:HD3	2.50	0.47
10:A:304:TDS:HBC1	2:B:85:PHE:CA	2.41	0.47
2:B:84:VAL:HG21	13:B:201:CLA:H41	1.96	0.47
3:P:177:THR:CA	3:P:227:ALA:HA	2.44	0.47
1:N:79:GLY:HA2	4:Q:41:TYR:CE1	2.49	0.47
3:P:77:MET:HG2	3:P:113:VAL:HG13	1.95	0.47
2:B:122:ASN:ND2	2:B:124:PHE:CD2	2.78	0.47
1:A:39:ILE:CD1	15:E:101:BCR:H312	2.41	0.47
8:H:22:VAL:C	8:H:24:TRP:N	2.67	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:C:211:ILE:HD11	3:C:231:LEU:HD23	1.95	0.47
3:C:223:GLN:O	3:C:224:ALA:HB3	2.14	0.47
1:A:188:THR:HG22	9:A:301:HEM:HBC2	1.95	0.47
3:C:198:SER:HA	3:C:208:VAL:HB	1.96	0.47
2:O:135:PHE:O	2:O:138:LEU:HB2	2.13	0.47
2:O:91:LEU:CD2	2:O:100:LEU:HD12	2.44	0.47
3:P:56:THR:HG22	3:P:122:GLU:HB3	1.95	0.47
4:D:55:THR:O	4:D:57:LYS:HG3	2.14	0.47
7:G:11:LEU:O	7:G:14:VAL:HG13	2.15	0.47
8:H:8:TRP:HA	8:H:11:LEU:HG	1.96	0.47
3:C:185:LYS:CD	3:C:195:TYR:HD2	2.26	0.47
2:O:95:LEU:HD23	2:O:95:LEU:H	1.77	0.47
1:N:21:VAL:CG1	1:N:22:THR:H	2.21	0.47
3:P:53:PRO:O	3:P:54:TYR:HB2	2.14	0.47
3:P:54:TYR:OH	3:P:58:LEU:HB2	2.14	0.47
2:B:27:TYR:CD1	2:B:27:TYR:N	2.83	0.47
3:C:283:GLN:C	3:C:285:ALA:N	2.67	0.47
3:P:50:VAL:O	3:P:50:VAL:HG12	2.14	0.47
2:O:41:PRO:O	2:O:45:MET:HG3	2.15	0.47
2:B:118:ASN:HB2	2:B:123:PRO:HG3	1.95	0.47
3:C:41:LEU:HD22	3:C:41:LEU:H	1.79	0.47
12:Q:1307:OPC:HAP2	12:Q:1307:OPC:HAL2	1.93	0.47
2:O:144:GLY:O	2:O:145:ILE:CD1	2.62	0.47
3:P:281:LYS:HD2	3:P:281:LYS:H	1.79	0.47
4:D:52:GLY:CA	4:D:85:LYS:NZ	2.77	0.47
2:O:127:PRO:O	2:O:129:ALA:N	2.41	0.47
1:A:170:ARG:HG3	1:A:172:GLY:H	1.79	0.47
3:C:26:HIS:ND1	3:C:154:ASN:ND2	2.62	0.47
1:A:62:VAL:HG11	1:A:177:GLN:HB3	1.96	0.47
1:N:47:GLN:HG2	9:N:301:HEM:HBB2	1.95	0.47
2:O:91:LEU:HD21	2:O:100:LEU:HD12	1.97	0.47
1:N:28:PRO:CG	2:O:32:TRP:CB	2.93	0.47
1:N:102:PHE:O	1:N:103:ARG:C	2.53	0.47
4:D:78:ARG:HB3	4:D:90:TYR:CE1	2.50	0.47
4:Q:134:ASP:HB2	4:Q:138:ARG:CB	2.43	0.47
1:A:202:HIS:O	1:A:206:ILE:HG13	2.15	0.47
5:E:25:ALA:C	5:E:27:LYS:H	2.18	0.47
8:H:5:VAL:O	8:H:9:VAL:CB	2.63	0.47
9:N:303:HEM:CBB	2:O:43:VAL:HG21	2.44	0.47
12:Q:1307:OPC:HAZ2	12:Q:1307:OPC:HBN2	1.97	0.47
1:N:186:ALA:HB2	10:N:1304:TDS:HAA3	1.97	0.47
2:O:148:ALA:O	2:O:149:LEU:O	2.32	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:N:141:ASP:C	2:O:64:GLU:CG	2.83	0.47
7:T:26:TYR:CA	7:T:29:TYR:CD1	2.90	0.47
7:T:25:ALA:O	7:T:29:TYR:CD2	2.68	0.47
4:Q:35:LEU:O	4:Q:35:LEU:HD12	2.15	0.47
1:A:103:ARG:NH2	9:A:302:HEM:HBD1	2.29	0.47
4:D:105:ASN:C	4:D:107:VAL:HG23	2.34	0.47
3:C:82:PHE:O	3:C:112:ASN:HB3	2.15	0.47
2:O:37:LEU:HD23	2:O:40:PHE:CE2	2.49	0.47
1:A:14:ILE:CG1	1:A:15:GLN:N	2.77	0.47
1:N:27:PRO:HB3	2:O:33:PRO:HD3	1.97	0.47
4:D:150:LEU:HD11	4:D:171:ARG:HD3	1.96	0.47
4:D:169:ASP:HB3	4:D:178:TRP:CZ2	2.50	0.47
11:A:305:PL9:HC71	11:A:305:PL9:H121	1.96	0.47
2:O:58:ASP:OD2	3:P:146:LYS:HE2	2.14	0.47
4:D:65:LYS:HB3	4:D:68:LYS:CD	2.42	0.47
2:B:144:GLY:O	2:B:145:ILE:CG1	2.63	0.47
2:B:148:ALA:O	2:B:152:ASP:N	2.48	0.47
3:C:13:PRO:O	3:C:21:VAL:HG22	2.14	0.47
1:N:120:SER:HA	1:N:123:ILE:HD12	1.97	0.47
4:D:93:VAL:CA	4:D:100:ARG:HB2	2.39	0.47
2:B:53:ALA:O	2:B:57:LEU:HG	2.14	0.47
1:A:95:LEU:HD13	1:A:96:MET:HE3	1.96	0.47
1:N:66:TYR:HE1	2:O:63:GLY:H	1.63	0.47
1:N:204:LEU:C	1:N:206:ILE:N	2.68	0.47
1:A:125:ALA:HB2	9:A:302:HEM:CMC	2.45	0.47
2:B:154:THR:HG22	2:B:155:LEU:N	2.30	0.47
1:A:91:SER:HB3	2:B:79:TRP:CZ3	2.50	0.47
3:C:1:TYR:O	3:C:4:TRP:HB2	2.15	0.47
4:D:84:LEU:HD23	4:D:84:LEU:C	2.35	0.47
3:C:49:VAL:HG12	3:C:51:LYS:HD2	1.96	0.47
3:P:41:LEU:HD22	3:P:41:LEU:H	1.79	0.47
1:A:141:ASP:O	1:A:145:TYR:N	2.42	0.47
1:A:144:GLY:O	1:A:148:VAL:HG12	2.14	0.47
1:A:15:GLN:CG	1:N:113:PRO:HA	2.44	0.47
1:A:146:TRP:HB2	2:B:75:ILE:CD1	2.43	0.47
4:Q:133:TYR:HA	4:Q:138:ARG:C	2.35	0.47
6:F:26:LEU:HD22	8:H:17:TRP:HE1	1.80	0.47
8:H:6:LEU:O	8:H:10:ALA:HB3	2.15	0.47
6:S:26:LEU:HD21	8:U:17:TRP:NE1	2.29	0.47
1:A:103:ARG:HH22	9:A:302:HEM:HBD1	1.80	0.47
1:A:91:SER:O	2:B:79:TRP:HH2	1.98	0.47
10:A:304:TDS:CBC	2:B:88:LEU:HD22	2.45	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:P:167:SER:O	3:P:172:PHE:CZ	2.68	0.47
3:P:198:SER:HA	3:P:208:VAL:HB	1.97	0.47
3:P:11:PRO:HA	3:P:107:LYS:NZ	2.29	0.47
3:P:275:LYS:HZ3	3:P:276:LYS:HZ1	1.60	0.46
5:R:23:ILE:HG23	5:R:24:PHE:HD2	1.76	0.46
4:D:99:ILE:HD13	4:D:153:ALA:O	2.15	0.46
4:Q:147:SER:HB3	4:Q:171:ARG:HH12	1.80	0.46
1:A:95:LEU:HD21	5:E:7:PHE:HB2	1.98	0.46
1:A:105:TYR:O	1:A:107:THR:N	2.48	0.46
1:A:15:GLN:HG3	1:N:113:PRO:HA	1.97	0.46
2:O:27:TYR:O	2:O:27:TYR:CD2	2.65	0.46
4:D:133:TYR:HE2	4:D:146:LEU:O	1.99	0.46
1:A:200:LEU:HD22	1:A:200:LEU:HA	1.76	0.46
1:A:204:LEU:C	1:A:206:ILE:N	2.68	0.46
4:D:50:VAL:O	4:D:164:PRO:HB2	2.15	0.46
5:E:18:ILE:HG22	5:E:22:ILE:CD1	2.45	0.46
3:P:231:LEU:CD1	3:P:233:ASN:H	2.28	0.46
3:P:79:PRO:HG2	3:P:82:PHE:CD1	2.50	0.46
4:Q:19:MET:C	4:Q:21:LEU:N	2.68	0.46
2:B:73:LEU:O	2:B:74:GLU:HB2	2.15	0.46
4:Q:63:ASN:O	4:Q:69:PHE:CE1	2.68	0.46
1:A:39:ILE:CD1	2:B:43:VAL:CG1	2.93	0.46
3:C:270:LEU:HG	3:C:274:LEU:CD1	2.45	0.46
12:Q:1307:OPC:HB2	12:Q:1307:OPC:CBX	2.38	0.46
2:O:145:ILE:HG23	2:O:149:LEU:HB2	1.98	0.46
4:D:156:GLN:HB2	4:D:161:VAL:CG2	2.45	0.46
3:P:199:ILE:HB	3:P:207:VAL:O	2.14	0.46
1:A:61:THR:HA	1:N:59:LYS:HZ2	1.80	0.46
11:A:305:PL9:C40	11:A:305:PL9:H512	2.42	0.46
3:C:277:LYS:HE3	7:G:31:ARG:CD	2.45	0.46
3:C:275:LYS:HE3	4:D:20:ASN:HB3	1.96	0.46
7:G:30:LYS:C	7:G:31:ARG:CZ	2.84	0.46
8:H:9:VAL:CG1	8:H:10:ALA:N	2.79	0.46
6:S:26:LEU:HB3	7:T:30:LYS:HE3	1.98	0.46
3:P:26:HIS:CB	3:P:154:ASN:HD21	2.29	0.46
3:C:250:GLN:NE2	3:C:251:ASP:H	2.12	0.46
2:O:60:ALA:C	2:O:61:MET:HG2	2.36	0.46
2:O:31:ALA:O	2:O:32:TRP:CB	2.54	0.46
4:D:168:THR:HA	4:D:176:PRO:HB3	1.97	0.46
4:Q:65:LYS:HB2	4:Q:68:LYS:HD3	1.98	0.46
3:C:268:ALA:HA	4:D:26:THR:OG1	2.15	0.46
7:G:14:VAL:HG22	7:G:15:PHE:N	2.28	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:N:95:LEU:HG	5:R:7:PHE:HB2	1.97	0.46
3:C:195:TYR:O	3:C:196:GLN:O	2.34	0.46
7:T:15:PHE:C	7:T:17:THR:H	2.18	0.46
3:C:93:GLU:CA	3:C:97:GLU:CB	2.87	0.46
1:A:105:TYR:C	1:A:107:THR:H	2.19	0.46
4:Q:53:GLY:O	4:Q:54:THR:OG1	2.27	0.46
4:D:15:ARG:NE	4:D:15:ARG:HA	2.31	0.46
2:B:42:VAL:CG1	7:G:28:GLN:HE22	2.28	0.46
5:E:9:ILE:HG21	6:F:9:ALA:C	2.34	0.46
5:E:5:ALA:HB2	6:F:5:MET:CG	2.46	0.46
3:C:231:LEU:CD1	3:C:231:LEU:H	2.29	0.46
4:D:151:CYS:SG	4:D:162:LEU:HD23	2.55	0.46
3:C:221:GLU:N	3:C:221:GLU:CD	2.68	0.46
6:F:3:GLU:HA	6:F:6:LEU:HD12	1.96	0.46
4:D:141:ARG:HG2	4:D:141:ARG:HH11	1.81	0.46
1:A:14:ILE:O	1:A:15:GLN:C	2.54	0.46
4:Q:12:ASP:O	4:Q:14:GLY:N	2.48	0.46
8:H:5:VAL:O	8:H:9:VAL:HG12	2.15	0.46
1:N:178:ALA:O	1:N:182:ARG:HG3	2.16	0.46
3:P:163:THR:OG1	3:P:165:GLU:HG2	2.16	0.46
8:U:5:VAL:O	8:U:9:VAL:HG12	2.16	0.46
1:A:188:THR:CG2	9:A:301:HEM:HBC2	2.46	0.46
3:C:181:THR:C	3:C:221:GLU:HB3	2.35	0.46
2:O:72:PRO:O	2:O:73:LEU:CB	2.62	0.46
1:N:27:PRO:HB3	2:O:31:ALA:HA	1.97	0.46
4:D:168:THR:HG22	4:D:169:ASP:N	2.30	0.46
4:Q:125:LYS:CD	4:Q:132:GLN:HG2	2.46	0.46
4:Q:165:TRP:HE1	4:Q:176:PRO:HB2	1.81	0.46
4:Q:64:VAL:HA	4:Q:69:PHE:CG	2.51	0.46
3:C:273:ILE:CG2	7:G:27:GLN:HB3	2.46	0.46
5:E:5:ALA:HB3	6:F:5:MET:HB2	1.98	0.46
7:T:15:PHE:C	7:T:17:THR:N	2.68	0.46
3:P:180:ILE:HG13	3:P:199:ILE:HA	1.97	0.46
3:C:47:LYS:HG3	3:C:128:VAL:HG13	1.97	0.46
5:R:18:ILE:HG22	5:R:22:ILE:CD1	2.46	0.46
5:R:9:ILE:CG2	6:S:12:SER:HB2	2.46	0.46
4:D:99:ILE:CD1	4:D:153:ALA:HB3	2.46	0.46
4:Q:117:TRP:HE1	4:Q:135:GLU:CB	2.29	0.46
4:Q:133:TYR:CA	4:Q:139:VAL:HA	2.38	0.46
4:Q:91:ILE:HD12	4:Q:160:ILE:CD1	2.38	0.46
3:C:265:VAL:O	3:C:269:GLN:HG3	2.14	0.46
7:G:28:GLN:C	7:G:30:LYS:N	2.68	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:P:167:SER:O	3:P:172:PHE:HZ	1.98	0.46
3:P:181:THR:C	3:P:221:GLU:HB3	2.36	0.46
2:O:21:GLY:C	2:O:24:HIS:HD2	2.19	0.46
3:C:59:GLN:OE1	3:C:67:LYS:HE3	2.16	0.46
4:D:36:TYR:O	4:D:40:LYS:HG2	2.16	0.46
4:D:57:LYS:H	4:D:82:GLN:HG2	1.81	0.46
4:Q:133:TYR:CD2	4:Q:148:LEU:HG	2.51	0.46
2:B:45:MET:HE3	4:D:27:VAL:HG13	1.97	0.46
6:F:7:TYR:CD1	6:F:7:TYR:C	2.89	0.46
2:O:87:ILE:HG13	2:O:143:LEU:HD13	1.98	0.46
1:N:141:ASP:HA	2:O:64:GLU:HG3	1.98	0.46
2:O:57:LEU:HD11	7:T:18:LEU:HB2	1.97	0.46
1:A:33:PHE:CD1	5:E:18:ILE:HG23	2.51	0.46
2:B:91:LEU:HD21	2:B:100:LEU:HD12	1.97	0.46
11:A:305:PL9:H18	12:B:307:OPC:HBU1	1.97	0.45
12:Q:1307:OPC:HBM1	12:Q:1307:OPC:HAR2	1.97	0.45
5:R:7:PHE:O	5:R:11:PHE:HD2	1.98	0.45
3:P:93:GLU:HG3	3:P:94:LEU:N	2.31	0.45
2:B:149:LEU:H	2:B:149:LEU:HG	1.26	0.45
2:B:104:VAL:HB	2:B:105:PRO:HD3	1.97	0.45
4:Q:22:LEU:HD23	4:Q:22:LEU:C	2.36	0.45
4:D:123:LYS:CB	4:D:134:ASP:HA	2.46	0.45
4:Q:115:VAL:HG13	4:Q:126:CYS:CB	2.44	0.45
3:C:266:MET:HE3	7:G:23:TYR:HB3	1.97	0.45
8:H:3:ILE:HG23	8:H:4:ASP:H	1.81	0.45
1:N:98:ILE:CD1	13:O:1201:CLA:HED3	2.45	0.45
2:O:102:ALA:O	2:O:105:PRO:HD2	2.16	0.45
2:O:57:LEU:HD21	3:P:258:MET:HE2	1.98	0.45
3:C:177:THR:CA	3:C:227:ALA:HA	2.45	0.45
3:P:1:TYR:HD2	3:P:119:LEU:CD2	2.25	0.45
3:C:163:THR:OG1	3:C:165:GLU:HG2	2.16	0.45
6:S:30:GLN:HB2	6:S:30:GLN:HE21	1.55	0.45
1:A:68:SER:O	1:A:71:TYR:HB3	2.15	0.45
1:A:15:GLN:HE21	1:A:15:GLN:HB2	1.57	0.45
3:P:282:VAL:HG22	4:Q:15:ARG:HE	1.82	0.45
4:D:168:THR:HG23	4:D:174:GLU:O	2.17	0.45
1:A:143:VAL:HA	2:B:75:ILE:CD1	2.46	0.45
4:Q:121:GLU:CD	4:Q:125:LYS:HE3	2.37	0.45
5:E:23:ILE:O	5:E:27:LYS:CB	2.60	0.45
7:G:10:VAL:O	7:G:14:VAL:HG13	2.17	0.45
2:O:79:TRP:CZ2	5:R:1:MET:CB	2.86	0.45
4:D:66:VAL:HG23	4:D:159:ASN:N	2.32	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:149:LEU:HD13	2:B:150:PRO:CD	2.43	0.45
1:A:106:LEU:HD12	5:E:18:ILE:HG13	1.99	0.45
3:P:3:PHE:HE1	3:P:117:GLY:CA	2.29	0.45
1:A:112:LYS:HG3	1:A:113:PRO:N	2.31	0.45
4:Q:100:ARG:HB3	4:Q:102:TYR:CE2	2.52	0.45
4:Q:175:LYS:HZ3	4:Q:175:LYS:HB3	1.82	0.45
1:A:23:SER:C	1:A:25:TYR:H	2.18	0.45
4:D:51:GLY:C	4:D:85:LYS:NZ	2.70	0.45
2:B:98:VAL:HA	2:B:101:MET:CE	2.46	0.45
1:A:142:GLN:HG2	2:B:65:PRO:O	2.16	0.45
2:B:26:TYR:O	2:B:27:TYR:HD1	1.98	0.45
2:O:40:PHE:N	2:O:41:PRO:HD2	2.31	0.45
2:O:28:GLY:O	2:O:29:GLU:OE2	2.35	0.45
4:D:108:CYS:O	4:D:110:HIS:N	2.50	0.45
4:D:123:LYS:NZ	4:D:140:ILE:CD1	2.80	0.45
4:Q:81:VAL:HB	4:Q:89:THR:HB	1.97	0.45
7:T:25:ALA:O	7:T:29:TYR:CG	2.70	0.45
2:B:83:PRO:O	2:B:87:ILE:HG13	2.15	0.45
8:U:9:VAL:HG22	8:U:13:VAL:HG23	1.97	0.45
3:C:52:ILE:HG21	3:C:154:ASN:O	2.16	0.45
1:N:209:GLN:HB2	2:O:27:TYR:HB2	1.98	0.45
4:D:116:PRO:HG3	4:D:126:CYS:HA	1.97	0.45
4:D:58:ASP:HB3	4:D:62:ASN:H	1.82	0.45
4:D:55:THR:O	4:D:82:GLN:HG3	2.16	0.45
4:D:98:ALA:O	4:D:100:ARG:N	2.50	0.45
4:Q:56:ALA:O	4:Q:57:LYS:O	2.34	0.45
6:S:7:TYR:CD1	6:S:7:TYR:C	2.89	0.45
2:B:77:PRO:HG3	2:B:85:PHE:CB	2.47	0.45
2:B:84:VAL:HG13	2:B:101:MET:HB2	1.99	0.45
3:P:219:VAL:HG12	3:P:220:SER:N	2.32	0.45
8:U:8:TRP:CE3	8:U:11:LEU:HB2	2.52	0.45
3:P:55:ASP:OD1	3:P:57:LYS:CG	2.64	0.45
4:Q:118:ASN:O	4:Q:122:ASN:HA	2.17	0.45
3:P:98:VAL:O	3:P:99:GLY:O	2.34	0.45
2:O:114:ILE:C	2:O:116:ASN:N	2.70	0.45
4:Q:12:ASP:HB2	4:Q:13:MET:CE	2.47	0.45
4:Q:16:ARG:HH11	4:Q:16:ARG:HG2	1.80	0.45
4:D:108:CYS:C	4:D:110:HIS:N	2.67	0.45
4:D:123:LYS:HB3	4:D:134:ASP:HA	1.99	0.45
4:Q:105:ASN:O	4:Q:149:ALA:HB3	2.16	0.45
1:A:26:VAL:HB	2:B:29:GLU:HB2	1.98	0.45
4:D:38:LEU:HG	12:D:306:OPC:OBH	2.17	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:O:86:GLN:NE2	2:O:145:ILE:HB	2.31	0.45
4:D:65:LYS:CB	4:D:68:LYS:HD3	2.41	0.45
1:N:202:HIS:O	1:N:206:ILE:HG13	2.16	0.45
3:P:28:ALA:HB3	3:P:238:GLY:O	2.17	0.45
1:A:106:LEU:HD21	2:B:133:PHE:CZ	2.52	0.45
3:P:185:LYS:HE3	3:P:217:LEU:HD21	1.98	0.45
3:P:190:TYR:HB3	3:P:191:GLY:H	1.44	0.45
8:U:7:GLY:O	8:U:11:LEU:CD1	2.64	0.45
3:C:229:GLU:O	3:C:230:ALA:C	2.55	0.45
3:C:3:PHE:CZ	3:C:119:LEU:HD11	2.49	0.45
3:C:197:VAL:HG12	3:C:198:SER:N	2.32	0.45
3:P:40:VAL:HG11	3:P:46:PHE:CD2	2.51	0.45
3:C:59:GLN:HB2	3:C:67:LYS:HE3	1.97	0.45
1:A:119:ILE:C	1:A:121:GLY:N	2.67	0.45
3:P:251:ASP:HB3	3:P:254:ARG:HD2	1.98	0.45
4:Q:19:MET:O	4:Q:21:LEU:N	2.50	0.45
4:D:124:PHE:CD2	4:D:148:LEU:HD12	2.52	0.45
4:D:99:ILE:CB	4:D:153:ALA:HB1	2.36	0.45
4:D:58:ASP:OD1	4:D:69:PHE:CE2	2.70	0.45
4:D:57:LYS:N	4:D:82:GLN:HG2	2.32	0.45
11:Q:1305:PL9:H412	11:Q:1305:PL9:H512	1.99	0.45
3:C:171:VAL:HG13	3:C:234:ASN:CB	2.45	0.45
8:U:24:TRP:CG	8:U:25:GLY:N	2.83	0.45
3:P:280:GLU:HG2	3:P:281:LYS:HZ1	1.79	0.45
2:B:148:ALA:C	2:B:149:LEU:HD12	2.36	0.45
3:P:4:TRP:CE3	3:P:4:TRP:N	2.85	0.45
1:A:17:LEU:HA	1:A:20:ASP:OD1	2.17	0.45
4:Q:19:MET:C	4:Q:21:LEU:H	2.19	0.45
5:R:30:LYS:NZ	5:R:30:LYS:CB	2.80	0.45
4:Q:156:GLN:CB	4:Q:161:VAL:HG11	2.47	0.45
4:Q:165:TRP:HA	4:Q:167:GLU:OE2	2.17	0.45
1:A:28:PRO:HD2	2:B:31:ALA:C	2.36	0.45
1:N:154:VAL:CG1	10:N:1304:TDS:HAX1	2.47	0.45
1:N:137:SER:O	1:N:140:TRP:CD1	2.64	0.45
7:T:14:VAL:HG22	7:T:15:PHE:N	2.32	0.45
3:C:1:TYR:HD2	3:C:119:LEU:CD2	2.29	0.45
4:Q:16:ARG:HG2	4:Q:16:ARG:NH1	2.32	0.45
4:D:64:VAL:HG21	4:D:81:VAL:CG1	2.47	0.45
1:A:114:ARG:NH1	1:A:209:GLN:N	2.55	0.45
2:B:53:ALA:HB2	3:C:262:ILE:HD11	1.99	0.45
3:C:218:ILE:HG23	3:C:232:THR:O	2.17	0.45
2:O:57:LEU:HD22	7:T:14:VAL:CB	2.47	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:O:57:LEU:HD22	7:T:14:VAL:HB	1.99	0.45
3:P:93:GLU:HG3	3:P:94:LEU:H	1.82	0.45
8:U:9:VAL:C	8:U:11:LEU:N	2.69	0.45
3:P:14:ARG:HG2	3:P:77:MET:CE	2.47	0.45
1:N:111:LYS:HB3	1:N:113:PRO:HD2	1.99	0.44
4:Q:88:PRO:HD3	4:Q:107:VAL:HG23	1.99	0.44
4:Q:117:TRP:HE1	4:Q:135:GLU:CA	2.25	0.44
1:A:114:ARG:HB3	1:A:208:LYS:HZ3	1.82	0.44
8:H:11:LEU:O	8:H:12:LEU:C	2.54	0.44
3:C:187:GLU:CG	3:C:188:ASP:H	2.28	0.44
1:N:116:LEU:HD11	1:N:208:LYS:HZ1	1.82	0.44
1:A:71:TYR:CE1	1:A:75:GLU:HG2	2.52	0.44
4:Q:59:LYS:HG2	4:Q:79:VAL:HG13	1.99	0.44
3:C:84:ILE:HD12	3:C:130:PRO:HD2	1.99	0.44
2:O:122:ASN:HD22	2:O:125:ARG:NH1	2.14	0.44
2:O:32:TRP:CZ3	2:O:36:LEU:HA	2.50	0.44
4:D:147:SER:HB2	4:D:177:TRP:CH2	2.45	0.44
4:Q:149:ALA:O	4:Q:150:LEU:HD22	2.17	0.44
4:Q:81:VAL:HG12	4:Q:82:GLN:N	2.32	0.44
4:Q:90:TYR:O	4:Q:91:ILE:CD1	2.65	0.44
3:C:41:LEU:N	3:C:41:LEU:HD22	2.32	0.44
8:H:11:LEU:C	8:H:15:PHE:CE1	2.91	0.44
10:N:1304:TDS:CBC	2:O:81:LEU:HB3	2.47	0.44
2:O:147:ALA:C	2:O:149:LEU:N	2.69	0.44
3:P:86:PRO:HB2	3:P:88:GLU:OE1	2.17	0.44
2:B:145:ILE:O	2:B:147:ALA:N	2.50	0.44
3:C:60:GLN:OE1	3:C:156:GLY:HA2	2.17	0.44
8:U:5:VAL:O	8:U:9:VAL:CB	2.64	0.44
1:N:115:GLU:N	1:N:115:GLU:OE1	2.46	0.44
4:Q:15:ARG:NE	4:Q:15:ARG:HA	2.32	0.44
1:A:29:HIS:CG	1:A:30:VAL:N	2.85	0.44
7:G:30:LYS:C	7:G:31:ARG:NE	2.71	0.44
3:P:218:ILE:HG23	3:P:232:THR:O	2.17	0.44
1:N:77:SER:O	1:N:78:PHE:HB2	2.17	0.44
1:N:58:TYR:CE2	1:N:60:PRO:HB3	2.53	0.44
1:N:29:HIS:HE2	1:N:209:GLN:NE2	2.15	0.44
4:D:55:THR:HG1	4:D:57:LYS:HZ2	1.60	0.44
4:Q:102:TYR:N	4:Q:153:ALA:HB2	2.32	0.44
2:B:38:TYR:HB3	3:C:276:LYS:HD3	1.99	0.44
3:C:277:LYS:CE	7:G:31:ARG:NE	2.81	0.44
3:C:185:LYS:HD3	3:C:195:TYR:CD2	2.47	0.44
3:C:48:ALA:HB3	3:C:129:PHE:HB2	1.98	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:C:154:ASN:OD1	3:C:155:ARG:N	2.49	0.44
1:A:52:PHE:HB2	1:N:189:PHE:CE2	2.52	0.44
2:B:86:GLN:HG2	2:B:143:LEU:CA	2.47	0.44
3:P:41:LEU:HD22	3:P:41:LEU:N	2.32	0.44
4:Q:161:VAL:C	4:Q:162:LEU:HD12	2.38	0.44
5:E:15:PHE:O	5:E:16:PHE:C	2.56	0.44
7:T:26:TYR:HA	7:T:29:TYR:HD1	1.71	0.44
3:P:120:PRO:CB	3:P:124:TYR:CD1	3.01	0.44
2:B:86:GLN:O	2:B:90:SER:HB2	2.18	0.44
1:N:77:SER:O	4:Q:41:TYR:HA	2.18	0.44
3:P:71:ASN:HB2	9:P:301:HEM:CGA	2.47	0.44
5:R:23:ILE:O	5:R:27:LYS:HG2	2.17	0.44
4:D:57:LYS:CG	4:D:82:GLN:HG2	2.48	0.44
4:Q:109:THR:C	4:Q:145:PRO:HG3	2.37	0.44
4:Q:68:LYS:N	4:Q:68:LYS:CD	2.78	0.44
1:A:114:ARG:NH1	1:A:209:GLN:O	2.51	0.44
4:D:22:LEU:O	4:D:26:THR:HG23	2.17	0.44
3:C:190:TYR:HB3	3:C:191:GLY:H	1.42	0.44
12:N:1306:OPC:HBG1	4:Q:39:VAL:HG13	1.99	0.44
1:A:189:PHE:O	1:A:192:PRO:HG2	2.17	0.44
3:C:45:VAL:HG22	3:C:85:ALA:CB	2.48	0.44
1:A:32:ILE:C	1:A:34:TYR:N	2.68	0.44
1:N:159:PRO:C	1:N:160:VAL:HG23	2.38	0.44
1:N:28:PRO:C	1:N:29:HIS:ND1	2.71	0.44
4:D:77:ASP:OD2	4:D:77:ASP:C	2.56	0.44
4:Q:103:GLY:O	4:Q:105:ASN:N	2.50	0.44
4:Q:123:LYS:CD	4:Q:140:ILE:HD12	2.33	0.44
4:Q:81:VAL:C	4:Q:82:GLN:HG2	2.38	0.44
2:O:146:GLY:O	2:O:149:LEU:CG	2.65	0.44
2:O:81:LEU:O	2:O:85:PHE:CB	2.66	0.44
2:O:57:LEU:HD13	7:T:15:PHE:CD2	2.53	0.44
7:T:28:GLN:C	7:T:30:LYS:H	2.21	0.44
1:A:105:TYR:C	1:A:107:THR:N	2.71	0.44
8:U:22:VAL:HG12	8:U:26:ARG:HE	1.82	0.44
4:D:16:ARG:O	4:D:18:PHE:N	2.48	0.44
2:O:25:ASN:OD1	2:O:26:TYR:HB2	2.17	0.44
2:O:96:LEU:O	2:O:100:LEU:CB	2.66	0.44
1:A:119:ILE:O	1:A:123:ILE:HG13	2.18	0.44
5:R:26:ILE:N	5:R:26:ILE:CD1	2.80	0.44
4:D:117:TRP:CG	4:D:118:ASN:N	2.85	0.44
4:D:177:TRP:CE3	4:D:178:TRP:HE3	2.36	0.44
2:B:74:GLU:O	2:B:75:ILE:HB	2.18	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:Q:78:ARG:HG2	4:Q:117:TRP:CE3	2.53	0.44
4:Q:124:PHE:HD2	4:Q:133:TYR:HB2	1.82	0.44
4:Q:83:GLY:HA3	4:Q:87:ASP:O	2.17	0.44
1:A:27:PRO:HB3	2:B:31:ALA:C	2.38	0.44
1:N:92:MET:O	1:N:96:MET:HG2	2.18	0.44
2:O:146:GLY:H	2:O:149:LEU:CD1	2.22	0.44
2:O:146:GLY:O	2:O:147:ALA:C	2.55	0.44
8:U:9:VAL:CG1	8:U:10:ALA:N	2.81	0.44
3:C:173:THR:HB	3:C:229:GLU:HA	1.99	0.44
3:P:61:VAL:HG23	3:P:61:VAL:O	2.17	0.44
4:D:59:LYS:HD2	4:D:79:VAL:HB	1.99	0.44
1:N:44:PHE:HD1	9:N:301:HEM:HBB1	1.82	0.44
3:C:54:TYR:OH	3:C:58:LEU:HB2	2.18	0.44
1:N:81:LEU:HD13	1:N:81:LEU:O	2.18	0.44
1:N:119:ILE:C	1:N:121:GLY:N	2.69	0.44
4:D:138:ARG:CZ	4:D:138:ARG:CB	2.96	0.44
2:B:122:ASN:HA	2:B:123:PRO:HD3	1.69	0.44
2:B:32:TRP:CZ3	2:B:36:LEU:HA	2.47	0.44
7:G:9:LEU:CG	7:G:10:VAL:H	2.31	0.44
7:T:11:LEU:CA	7:T:14:VAL:HG13	2.47	0.44
3:P:233:ASN:HD22	3:P:234:ASN:H	1.66	0.44
1:A:159:PRO:C	1:A:161:VAL:N	2.68	0.44
4:D:152:HIS:N	4:D:152:HIS:CD2	2.86	0.44
5:R:5:ALA:HB2	6:S:5:MET:CB	2.48	0.43
4:D:100:ARG:HD3	4:D:101:ASP:N	2.33	0.43
4:D:118:ASN:HB3	4:D:121:GLU:H	1.83	0.43
4:D:129:HIS:O	4:D:143:PRO:CG	2.65	0.43
4:Q:163:THR:O	4:Q:164:PRO:C	2.57	0.43
4:Q:64:VAL:HB	4:Q:81:VAL:HG23	1.98	0.43
1:A:28:PRO:CD	2:B:32:TRP:HB2	2.48	0.43
3:C:279:VAL:O	3:C:279:VAL:HG12	2.16	0.43
2:O:145:ILE:HG22	2:O:146:GLY:N	2.33	0.43
3:C:231:LEU:CD1	3:C:231:LEU:N	2.80	0.43
3:P:262:ILE:HG13	7:T:21:LEU:HG	2.00	0.43
1:N:200:LEU:HA	1:N:200:LEU:HD22	1.76	0.43
2:B:149:LEU:O	2:B:150:PRO:C	2.56	0.43
2:B:151:LEU:HA	2:B:153:LYS:CD	2.42	0.43
8:U:11:LEU:O	8:U:12:LEU:C	2.56	0.43
3:C:201:THR:HG22	3:C:201:THR:O	2.18	0.43
3:P:48:ALA:HB3	3:P:129:PHE:HB2	1.99	0.43
5:R:27:LYS:HD3	5:R:27:LYS:HA	1.71	0.43
4:D:99:ILE:HD12	4:D:101:ASP:OD1	2.17	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:Q:115:VAL:CG2	4:Q:126:CYS:HB2	2.47	0.43
4:Q:94:GLU:N	4:Q:94:GLU:OE1	2.52	0.43
3:C:275:LYS:O	3:C:276:LYS:O	2.36	0.43
5:R:2:ILE:HG23	5:R:3:LEU:H	1.83	0.43
3:C:231:LEU:HD13	3:C:232:THR:N	2.26	0.43
1:N:140:TRP:O	2:O:65:PRO:CG	2.65	0.43
10:A:304:TDS:HBC3	2:B:88:LEU:HD22	2.00	0.43
3:C:26:HIS:CB	3:C:154:ASN:HD21	2.31	0.43
1:A:111:LYS:CG	1:A:112:LYS:N	2.73	0.43
3:P:45:VAL:HG22	3:P:85:ALA:CB	2.47	0.43
3:P:29:ALA:O	3:P:30:LYS:HG2	2.18	0.43
2:B:99:LEU:HD13	2:B:99:LEU:C	2.39	0.43
1:N:103:ARG:NE	1:N:107:THR:HG21	2.19	0.43
1:N:119:ILE:O	1:N:123:ILE:HG13	2.18	0.43
2:B:73:LEU:HD22	2:B:73:LEU:N	2.33	0.43
12:D:306:OPC:HBX1	1:N:201:LEU:HD13	1.99	0.43
8:H:21:MET:O	8:H:24:TRP:HE3	2.01	0.43
5:E:9:ILE:HB	6:F:9:ALA:HB1	2.00	0.43
2:B:149:LEU:CB	2:B:150:PRO:CD	2.95	0.43
3:P:180:ILE:HG23	3:P:220:SER:C	2.38	0.43
2:O:67:ASN:HD22	3:P:16:PRO:HB3	1.80	0.43
6:F:10:LEU:HD23	6:F:10:LEU:N	2.32	0.43
1:N:30:VAL:HG23	1:N:30:VAL:O	2.17	0.43
4:D:78:ARG:HH21	4:D:92:VAL:HG11	1.83	0.43
4:Q:91:ILE:HG22	4:Q:91:ILE:O	2.18	0.43
4:Q:76:GLY:CA	4:Q:93:VAL:O	2.62	0.43
1:A:30:VAL:O	1:A:30:VAL:CG2	2.66	0.43
12:B:307:OPC:CAL	12:B:307:OPC:HAP1	2.45	0.43
2:B:24:HIS:O	2:B:25:ASN:HB3	2.18	0.43
2:B:67:ASN:HA	2:B:68:PRO:HD2	1.80	0.43
9:A:301:HEM:HBD1	9:A:301:HEM:CHA	2.47	0.43
3:P:1:TYR:CD2	3:P:119:LEU:HD21	2.48	0.43
3:C:220:SER:HB2	3:C:221:GLU:CD	2.39	0.43
3:P:213:ALA:C	3:P:215:PRO:HD2	2.38	0.43
1:N:118:TRP:CZ3	2:O:108:LEU:C	2.91	0.43
4:Q:16:ARG:NH2	4:Q:18:PHE:HE1	2.16	0.43
4:Q:19:MET:HG3	4:Q:20:ASN:N	2.34	0.43
5:R:18:ILE:HA	5:R:22:ILE:HG13	2.01	0.43
4:D:81:VAL:HG21	4:D:91:ILE:CD1	2.43	0.43
4:D:57:LYS:H	4:D:82:GLN:HB2	1.84	0.43
4:Q:91:ILE:O	4:Q:92:VAL:O	2.35	0.43
8:H:8:TRP:O	8:H:11:LEU:HB2	2.17	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:E:7:PHE:CD1	5:E:7:PHE:C	2.92	0.43
5:E:9:ILE:CG1	6:F:13:PHE:HB2	2.47	0.43
2:O:136:GLY:HA3	13:O:1201:CLA:C4B	2.48	0.43
2:O:57:LEU:HD12	7:T:18:LEU:HD12	2.00	0.43
2:B:148:ALA:O	2:B:149:LEU:C	2.57	0.43
2:O:94:LYS:HB3	2:O:95:LEU:H	1.66	0.43
4:D:105:ASN:CB	4:D:107:VAL:HG23	2.49	0.43
4:Q:120:ALA:C	4:Q:122:ASN:N	2.71	0.43
4:D:73:HIS:ND1	4:D:77:ASP:OD1	2.52	0.43
4:Q:152:HIS:HB2	4:Q:162:LEU:HB3	2.01	0.43
4:Q:82:GLN:O	4:Q:87:ASP:O	2.37	0.43
3:C:255:VAL:O	3:C:258:MET:HG3	2.18	0.43
5:E:27:LYS:HD3	5:E:27:LYS:HA	1.75	0.43
2:B:25:ASN:ND2	2:B:25:ASN:N	2.66	0.43
12:Q:1307:OPC:HBE1	12:Q:1307:OPC:HBX2	2.01	0.43
4:D:68:LYS:H	4:D:68:LYS:CD	2.32	0.43
3:P:231:LEU:O	3:P:232:THR:CG2	2.58	0.43
1:A:142:GLN:HE21	2:B:66:ALA:CA	2.22	0.43
1:N:22:THR:O	1:N:25:TYR:CE2	2.71	0.43
3:C:14:ARG:HB3	3:C:20:ILE:HD13	2.01	0.43
3:P:41:LEU:HB3	3:P:42:PRO:HD2	2.00	0.43
1:N:34:TYR:CD2	1:N:34:TYR:N	2.87	0.43
4:D:78:ARG:HG3	4:D:117:TRP:CD1	2.54	0.43
4:D:118:ASN:CB	4:D:122:ASN:H	2.31	0.43
4:D:118:ASN:HB2	4:D:122:ASN:N	2.34	0.43
4:Q:92:VAL:HG23	4:Q:100:ARG:H	1.83	0.43
4:Q:115:VAL:CG1	4:Q:126:CYS:HB2	2.46	0.43
4:Q:134:ASP:N	4:Q:138:ARG:O	2.51	0.43
4:Q:155:VAL:HG13	4:Q:159:ASN:O	2.18	0.43
1:A:25:TYR:O	1:A:26:VAL:HG13	2.19	0.43
3:C:262:ILE:HG21	7:G:17:THR:HG22	1.99	0.43
2:B:24:HIS:O	2:B:25:ASN:CB	2.66	0.43
5:E:9:ILE:CG2	6:F:12:SER:HB2	2.47	0.43
2:O:141:ILE:CD1	5:R:11:PHE:CZ	3.01	0.43
3:P:262:ILE:HG22	3:P:263:CYS:N	2.34	0.43
1:A:106:LEU:HD12	5:E:18:ILE:CG1	2.48	0.43
3:P:211:ILE:HD11	3:P:231:LEU:HD23	1.99	0.43
4:D:16:ARG:NH2	4:D:18:PHE:HE1	2.17	0.43
2:O:22:MET:O	2:O:23:GLY:C	2.56	0.43
1:N:109:GLY:C	1:N:111:LYS:N	2.72	0.43
1:N:35:CYS:O	1:N:39:ILE:HG12	2.18	0.43
2:O:122:ASN:ND2	5:R:26:ILE:CG2	2.81	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:D:165:TRP:CZ2	4:D:178:TRP:NE1	2.87	0.43
4:Q:154:THR:C	4:Q:155:VAL:HG23	2.38	0.43
2:B:45:MET:HE1	4:D:27:VAL:HA	2.01	0.43
1:A:91:SER:OG	2:B:79:TRP:HZ3	2.02	0.43
4:D:12:ASP:HB2	4:D:13:MET:CE	2.49	0.43
2:B:94:LYS:HE2	2:B:94:LYS:HB2	1.82	0.43
4:Q:27:VAL:C	4:Q:29:GLY:N	2.71	0.43
4:Q:36:TYR:O	4:Q:40:LYS:HG2	2.18	0.43
1:N:110:PHE:HB3	1:N:118:TRP:CD1	2.54	0.43
1:N:118:TRP:HZ3	2:O:108:LEU:O	2.02	0.43
3:P:275:LYS:O	3:P:276:LYS:O	2.36	0.43
4:D:117:TRP:CZ2	4:D:122:ASN:HA	2.54	0.43
4:Q:78:ARG:HH12	4:Q:100:ARG:NH2	2.12	0.43
2:B:126:ARG:HG2	2:B:128:VAL:HG23	2.01	0.43
1:A:27:PRO:HD2	2:B:29:GLU:HB2	2.00	0.43
11:A:305:PL9:C10	12:D:306:OPC:HBF2	2.45	0.43
8:H:22:VAL:O	8:H:26:ARG:HG3	2.19	0.43
3:C:193:VAL:HG12	3:C:195:TYR:CE1	2.53	0.43
4:D:50:VAL:O	4:D:164:PRO:CB	2.66	0.43
1:A:44:PHE:CD1	9:A:301:HEM:HBB1	2.37	0.43
3:C:250:GLN:NE2	3:C:251:ASP:N	2.67	0.43
7:T:33:ASN:ND2	7:T:34:GLU:H	2.17	0.43
4:D:73:HIS:CE1	4:D:77:ASP:OD1	2.72	0.43
4:Q:133:TYR:HA	4:Q:139:VAL:N	2.34	0.43
4:Q:73:HIS:CE1	4:Q:77:ASP:HB2	2.53	0.43
7:T:28:GLN:O	7:T:29:TYR:C	2.57	0.43
3:P:88:GLU:CD	3:P:88:GLU:H	2.21	0.43
2:O:126:ARG:NH1	2:O:129:ALA:HB2	2.33	0.43
1:A:189:PHE:O	1:A:193:TRP:CE3	2.65	0.43
3:C:176:ALA:HB1	3:C:201:THR:HG21	2.00	0.43
2:O:116:ASN:HD22	2:O:116:ASN:C	2.21	0.42
2:O:18:LEU:HD22	2:O:31:ALA:CB	2.49	0.42
4:D:55:THR:O	4:D:55:THR:OG1	2.32	0.42
4:D:78:ARG:NE	4:D:92:VAL:CG1	2.78	0.42
4:Q:146:LEU:O	4:Q:147:SER:O	2.37	0.42
4:Q:175:LYS:NZ	4:Q:175:LYS:CB	2.82	0.42
3:C:270:LEU:HG	3:C:274:LEU:HD11	2.00	0.42
5:E:16:PHE:CD2	15:E:101:BCR:H361	2.54	0.42
6:F:22:LEU:O	6:F:26:LEU:HB2	2.18	0.42
4:D:66:VAL:O	4:D:70:LEU:HB2	2.18	0.42
1:A:186:ALA:HA	1:A:190:VAL:HB	1.99	0.42
1:A:91:SER:CB	2:B:79:TRP:CZ3	3.02	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:142:TRP:HD1	2:B:155:LEU:C	2.20	0.42
3:P:161:TYR:C	3:P:163:THR:H	2.21	0.42
9:A:301:HEM:HMA1	9:A:301:HEM:HAA2	1.87	0.42
2:B:95:LEU:N	2:B:95:LEU:CD2	2.71	0.42
1:N:131:PHE:CD1	9:N:301:HEM:HAB	2.53	0.42
1:A:14:ILE:HD13	1:A:15:GLN:CA	2.49	0.42
1:N:112:LYS:HG2	1:N:112:LYS:H	1.43	0.42
4:D:123:LYS:HZ2	4:D:140:ILE:CD1	2.32	0.42
4:D:168:THR:OG1	4:D:176:PRO:HD3	2.18	0.42
1:N:154:VAL:N	1:N:155:PRO:HD2	2.34	0.42
1:N:186:ALA:HA	1:N:190:VAL:HB	2.00	0.42
1:N:95:LEU:HD21	5:R:7:PHE:HB2	2.01	0.42
3:C:117:GLY:N	3:C:118:PRO:HD2	2.33	0.42
3:P:154:ASN:OD1	3:P:155:ARG:N	2.51	0.42
1:N:68:SER:O	1:N:71:TYR:HB3	2.19	0.42
3:C:54:TYR:CD2	3:C:70:LEU:HD13	2.52	0.42
1:N:13:GLU:HB3	1:N:14:ILE:H	1.53	0.42
1:A:141:ASP:OD2	1:A:144:GLY:N	2.49	0.42
3:P:27:LEU:HG	3:P:27:LEU:H	1.26	0.42
4:D:166:THR:HA	4:D:179:VAL:HG22	2.01	0.42
4:Q:138:ARG:CZ	4:Q:138:ARG:HA	2.48	0.42
2:B:145:ILE:HG22	2:B:145:ILE:O	2.20	0.42
3:P:197:VAL:HG12	3:P:198:SER:N	2.33	0.42
3:C:14:ARG:HG2	3:C:77:MET:CE	2.49	0.42
1:N:34:TYR:CD2	1:N:103:ARG:HD3	2.54	0.42
4:Q:155:VAL:CG1	4:Q:156:GLN:N	2.82	0.42
4:Q:91:ILE:O	4:Q:99:ILE:HB	2.20	0.42
6:F:30:GLN:HE21	6:F:30:GLN:HB2	1.56	0.42
5:E:18:ILE:HB	5:E:22:ILE:HD12	2.01	0.42
8:U:3:ILE:HG23	8:U:4:ASP:H	1.85	0.42
1:A:135:GLY:HA2	1:A:138:LEU:HG	2.00	0.42
3:P:153:ALA:O	3:P:240:PHE:CD1	2.72	0.42
3:P:245:THR:HB	6:S:1:MET:HE1	2.01	0.42
4:Q:73:HIS:HB3	4:Q:93:VAL:CB	2.50	0.42
4:Q:73:HIS:HD2	4:Q:92:VAL:C	2.23	0.42
2:O:86:GLN:HG2	2:O:143:LEU:CA	2.48	0.42
4:D:155:VAL:CG1	4:D:156:GLN:N	2.83	0.42
2:O:64:GLU:CG	2:O:65:PRO:CD	2.86	0.42
6:S:18:VAL:O	6:S:22:LEU:HG	2.18	0.42
4:D:52:GLY:CA	4:D:85:LYS:HZ3	2.33	0.42
3:C:11:PRO:HA	3:C:107:LYS:HZ3	1.83	0.42
1:A:21:VAL:HG12	1:A:21:VAL:O	2.18	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:N:211:ILE:HB	1:N:212:SER:H	1.62	0.42
3:P:275:LYS:HG3	4:Q:20:ASN:HB3	2.00	0.42
4:Q:65:LYS:HD3	4:Q:68:LYS:CG	2.48	0.42
4:Q:76:GLY:N	4:Q:93:VAL:CG1	2.72	0.42
7:G:17:THR:O	7:G:21:LEU:HB3	2.19	0.42
1:N:95:LEU:CD2	5:R:7:PHE:HB2	2.50	0.42
3:C:92:GLU:N	3:C:94:LEU:HD23	2.35	0.42
1:N:83:ARG:HD2	9:N:301:HEM:O2D	2.19	0.42
1:A:189:PHE:CD1	1:A:193:TRP:CZ3	3.08	0.42
1:N:149:LYS:HD2	1:N:175:VAL:HG21	2.01	0.42
3:P:201:THR:C	3:P:203:SER:N	2.73	0.42
1:A:182:ARG:HH21	1:N:57:TYR:HE2	1.67	0.42
4:Q:36:TYR:N	4:Q:37:PRO:HD2	2.34	0.42
7:T:33:ASN:ND2	7:T:34:GLU:N	2.68	0.42
4:Q:155:VAL:HG22	4:Q:160:ILE:HA	2.02	0.42
3:C:273:ILE:CG2	7:G:31:ARG:HH12	2.32	0.42
5:E:23:ILE:CG2	5:E:24:PHE:H	2.32	0.42
7:G:23:TYR:O	7:G:24:ALA:C	2.58	0.42
2:O:80:TYR:HB3	13:O:1201:CLA:CED	2.38	0.42
3:C:172:PHE:H	3:C:231:LEU:CG	2.08	0.42
3:P:161:TYR:C	3:P:163:THR:N	2.73	0.42
1:A:173:SER:HB2	1:A:174:SER:H	1.60	0.42
3:C:177:THR:CB	3:C:227:ALA:HA	2.49	0.42
1:A:138:LEU:C	1:A:140:TRP:H	2.23	0.42
1:N:59:LYS:O	1:N:65:ALA:HA	2.19	0.42
1:N:81:LEU:O	1:N:85:ILE:HG13	2.20	0.42
2:B:135:PHE:O	2:B:139:VAL:HG23	2.20	0.42
1:A:141:ASP:O	1:A:145:TYR:HB3	2.20	0.42
7:T:33:ASN:CG	7:T:34:GLU:H	2.23	0.42
1:N:28:PRO:HD2	2:O:32:TRP:C	2.40	0.42
2:O:122:ASN:ND2	2:O:125:ARG:CZ	2.74	0.42
3:P:275:LYS:HD3	3:P:276:LYS:N	2.34	0.42
4:D:138:ARG:H	4:D:138:ARG:HG2	1.45	0.42
4:Q:165:TRP:HE1	4:Q:176:PRO:CB	2.33	0.42
2:B:38:TYR:HA	2:B:38:TYR:HD2	1.62	0.42
3:C:277:LYS:O	3:C:280:GLU:HB3	2.20	0.42
3:C:280:GLU:O	3:C:284:ALA:HB2	2.20	0.42
7:G:22:PHE:O	7:G:26:TYR:CE1	2.73	0.42
2:O:87:ILE:O	2:O:88:LEU:C	2.58	0.42
3:C:88:GLU:CD	3:C:88:GLU:H	2.23	0.42
3:P:92:GLU:N	3:P:94:LEU:HD23	2.33	0.42
1:A:88:TRP:HZ2	2:B:58:ASP:CB	2.31	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:P:185:LYS:CG	3:P:195:TYR:HB3	2.32	0.42
3:P:171:VAL:HB	3:P:231:LEU:CD1	2.49	0.42
2:O:94:LYS:HE2	2:O:94:LYS:HB2	1.84	0.42
3:P:54:TYR:CD2	3:P:70:LEU:HD13	2.48	0.42
6:S:34:LYS:HE2	6:S:34:LYS:N	2.35	0.42
2:B:91:LEU:CD2	2:B:100:LEU:HD12	2.50	0.42
7:T:27:GLN:CA	7:T:27:GLN:NE2	2.83	0.42
3:P:84:ILE:HG13	3:P:103:PHE:HD1	1.84	0.42
2:O:29:GLU:OE2	2:O:29:GLU:HA	2.19	0.42
4:D:147:SER:CB	4:D:171:ARG:HH22	2.33	0.42
4:Q:88:PRO:HD2	4:Q:107:VAL:HG23	2.02	0.42
4:Q:55:THR:O	4:Q:56:ALA:CB	2.67	0.42
4:Q:73:HIS:CD2	4:Q:77:ASP:CB	2.92	0.42
4:Q:87:ASP:HB3	4:Q:88:PRO:CD	2.43	0.42
1:A:29:HIS:CD2	1:A:30:VAL:CG2	3.00	0.42
2:B:37:LEU:O	2:B:38:TYR:CD2	2.73	0.42
7:G:28:GLN:O	7:G:31:ARG:HD2	2.20	0.42
11:Q:1305:PL9:H512	11:Q:1305:PL9:C41	2.50	0.42
2:B:77:PRO:O	2:B:78:GLU:C	2.57	0.42
3:P:52:ILE:HD12	3:P:127:ILE:HD12	2.02	0.42
1:N:116:LEU:H	1:N:116:LEU:CD1	2.16	0.42
1:A:191:LEU:N	1:A:192:PRO:HD2	2.35	0.42
3:P:248:VAL:HG12	3:P:250:GLN:HG2	2.01	0.42
3:P:50:VAL:HG21	3:P:76:LEU:HD22	2.02	0.42
3:C:22:CYS:O	3:C:24:ASN:N	2.53	0.42
3:P:268:ALA:HA	4:Q:26:THR:OG1	2.20	0.42
1:N:114:ARG:HH21	1:N:209:GLN:CA	2.33	0.42
4:D:131:SER:HB3	4:D:144:ALA:HA	2.01	0.42
4:D:165:TRP:CE2	4:D:167:GLU:O	2.72	0.42
4:D:177:TRP:CZ2	4:D:178:TRP:HZ3	2.38	0.42
3:C:248:VAL:HG12	3:C:250:GLN:HG2	2.01	0.42
3:P:59:GLN:HB2	3:P:67:LYS:CE	2.50	0.42
1:N:110:PHE:HB3	1:N:118:TRP:HB2	2.02	0.41
4:D:108:CYS:HB3	4:D:115:VAL:HG22	2.01	0.41
4:Q:64:VAL:O	4:Q:81:VAL:HG22	2.19	0.41
1:N:95:LEU:HD11	5:R:10:VAL:CG1	2.36	0.41
10:A:304:TDS:HBC2	2:B:85:PHE:HD2	1.84	0.41
3:P:182:LYS:CG	3:P:198:SER:HB2	2.46	0.41
3:P:154:ASN:O	3:P:155:ARG:HB3	2.19	0.41
4:D:16:ARG:HH11	4:D:16:ARG:HG2	1.85	0.41
3:P:59:GLN:HB2	3:P:67:LYS:HE3	2.01	0.41
2:O:124:PHE:HB2	5:R:23:ILE:HA	2.01	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:R:23:ILE:O	5:R:27:LYS:CG	2.68	0.41
4:D:118:ASN:N	4:D:123:LYS:O	2.53	0.41
4:Q:100:ARG:NH2	4:Q:135:GLU:O	2.50	0.41
4:D:15:ARG:HH21	4:D:17:GLN:HG2	1.82	0.41
2:O:104:VAL:HB	2:O:105:PRO:HD3	2.02	0.41
3:C:231:LEU:HD13	3:C:233:ASN:H	1.85	0.41
4:D:156:GLN:OE1	4:D:161:VAL:HG11	2.20	0.41
1:A:103:ARG:O	1:A:107:THR:HG22	2.20	0.41
2:B:79:TRP:HA	2:B:82:TYR:CZ	2.54	0.41
3:C:183:ILE:HG13	3:C:183:ILE:H	1.72	0.41
3:P:201:THR:O	3:P:201:THR:HG22	2.20	0.41
1:N:118:TRP:HZ3	2:O:108:LEU:C	2.23	0.41
4:D:55:THR:O	4:D:56:ALA:C	2.58	0.41
4:Q:109:THR:HG21	4:Q:177:TRP:CH2	2.54	0.41
4:Q:77:ASP:N	4:Q:77:ASP:OD1	2.53	0.41
5:E:26:ILE:CD1	5:E:26:ILE:N	2.83	0.41
11:Q:1305:PL9:H151	11:Q:1305:PL9:H171	1.68	0.41
11:Q:1305:PL9:H302	12:Q:1307:OPC:HBB2	2.02	0.41
3:C:93:GLU:HG3	3:C:94:LEU:N	2.35	0.41
3:P:185:LYS:CD	3:P:217:LEU:HD11	2.49	0.41
3:P:231:LEU:HD21	3:P:233:ASN:O	2.20	0.41
3:C:59:GLN:HB2	3:C:67:LYS:CE	2.51	0.41
1:A:77:SER:O	1:A:78:PHE:CB	2.67	0.41
1:A:50:THR:O	1:A:54:MET:HG3	2.21	0.41
1:A:146:TRP:CH2	2:B:71:THR:OG1	2.59	0.41
9:A:303:HEM:HBD2	9:A:303:HEM:HHA	2.02	0.41
2:B:22:MET:N	2:B:24:HIS:CE1	2.88	0.41
10:N:1304:TDS:CBB	2:O:81:LEU:HD13	2.47	0.41
5:R:7:PHE:C	5:R:7:PHE:CD1	2.94	0.41
2:B:150:PRO:O	2:B:153:LYS:CD	2.68	0.41
3:P:221:GLU:CD	3:P:221:GLU:N	2.74	0.41
3:P:231:LEU:HD22	3:P:232:THR:H	1.82	0.41
1:A:47:GLN:NE2	1:A:89:SER:HB3	2.29	0.41
3:P:84:ILE:HG23	3:P:130:PRO:HG2	2.01	0.41
1:A:14:ILE:HA	1:A:17:LEU:CD1	2.50	0.41
4:Q:15:ARG:HH21	4:Q:17:GLN:HG2	1.84	0.41
4:D:93:VAL:HG22	4:D:98:ALA:O	2.21	0.41
2:B:71:THR:CB	2:B:72:PRO:CD	2.80	0.41
8:H:25:GLY:O	8:H:29:LEU:HG	2.20	0.41
2:O:78:GLU:HG2	2:O:80:TYR:HD1	1.85	0.41
3:P:93:GLU:CA	3:P:97:GLU:CB	2.94	0.41
1:A:104:VAL:HG11	9:A:302:HEM:HMD2	2.03	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:P:229:GLU:O	3:P:230:ALA:C	2.58	0.41
3:C:118:PRO:O	3:C:120:PRO:HD3	2.20	0.41
6:S:34:LYS:HD2	6:S:35:GLU:N	2.35	0.41
1:N:158:ILE:HD11	2:O:98:VAL:HG11	2.03	0.41
2:O:155:LEU:HG	2:O:155:LEU:O	2.21	0.41
3:C:13:PRO:HB3	3:C:106:TYR:CE1	2.54	0.41
4:Q:50:VAL:O	4:Q:50:VAL:HG12	2.21	0.41
2:B:110:LEU:O	2:B:113:PHE:HB2	2.20	0.41
1:N:34:TYR:N	1:N:34:TYR:HD2	2.19	0.41
5:R:9:ILE:HG21	6:S:9:ALA:C	2.33	0.41
4:D:78:ARG:NH1	4:D:124:PHE:CZ	2.88	0.41
4:Q:139:VAL:O	4:Q:139:VAL:HG13	2.21	0.41
2:O:81:LEU:O	2:O:85:PHE:HB2	2.20	0.41
1:N:140:TRP:CZ2	1:N:145:TYR:HD2	2.39	0.41
1:N:142:GLN:HG2	2:O:64:GLU:HB3	2.03	0.41
7:T:11:LEU:H	7:T:11:LEU:CD2	2.24	0.41
1:A:105:TYR:CE1	13:B:201:CLA:HAB	2.55	0.41
1:A:156:GLU:HG2	1:A:163:VAL:HG22	2.02	0.41
3:P:52:ILE:CG1	3:P:153:ALA:HB1	2.44	0.41
3:C:159:GLN:HG3	3:C:170:ASN:HB3	2.02	0.41
4:D:167:GLU:N	4:D:167:GLU:OE2	2.53	0.41
4:Q:78:ARG:HH21	4:Q:135:GLU:HB2	1.85	0.41
12:B:307:OPC:HBS	12:B:307:OPC:HBC1	2.02	0.41
8:H:15:PHE:O	8:H:19:ILE:HG13	2.21	0.41
2:B:89:ARG:NH2	2:B:147:ALA:O	2.51	0.41
1:A:33:PHE:HZ	5:E:17:GLY:HA3	1.85	0.41
3:P:187:GLU:CG	3:P:188:ASP:H	2.32	0.41
3:C:177:THR:OG1	3:C:227:ALA:HA	2.21	0.41
1:A:47:GLN:OE1	1:A:86:HIS:O	2.39	0.41
4:D:39:VAL:O	4:D:43:ILE:HD13	2.20	0.41
7:T:33:ASN:HD22	7:T:33:ASN:H	1.68	0.41
2:O:34:ASN:OD1	2:O:35:ASP:N	2.54	0.41
3:P:73:GLY:O	3:P:74:ALA:HB2	2.21	0.41
5:R:22:ILE:HA	5:R:25:ALA:HB2	2.02	0.41
4:D:126:CYS:HB2	4:D:131:SER:HB2	2.02	0.41
4:Q:137:GLY:C	4:Q:171:ARG:NH1	2.73	0.41
4:Q:168:THR:HG23	4:Q:175:LYS:CA	2.50	0.41
1:A:29:HIS:CE1	1:A:210:GLY:O	2.74	0.41
2:B:40:PHE:N	2:B:41:PRO:CD	2.84	0.41
7:G:11:LEU:H	7:G:11:LEU:CD2	2.23	0.41
3:C:86:PRO:HB2	3:C:88:GLU:OE1	2.21	0.41
3:P:161:TYR:CE1	3:P:167:SER:CA	2.96	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:P:180:ILE:HG12	3:P:181:THR:N	2.36	0.41
3:P:192:ASN:OD1	3:P:193:VAL:N	2.54	0.41
3:P:183:ILE:HG13	3:P:183:ILE:H	1.72	0.41
3:P:10:PRO:N	3:P:11:PRO:HD2	2.36	0.41
1:N:148:VAL:HA	1:N:151:VAL:HG22	2.03	0.41
5:E:30:LYS:HB3	5:E:30:LYS:HZ3	1.86	0.41
1:A:120:SER:HA	1:A:123:ILE:HD12	2.03	0.41
3:P:264:LEU:HD13	4:Q:30:VAL:HA	2.02	0.41
1:N:39:ILE:CD1	15:R:1101:BCR:H312	2.50	0.41
5:R:23:ILE:CG2	5:R:24:PHE:H	2.34	0.41
4:D:81:VAL:CG1	4:D:82:GLN:H	2.20	0.41
4:D:138:ARG:NH2	4:D:147:SER:CB	2.83	0.41
4:Q:78:ARG:CD	4:Q:78:ARG:N	2.82	0.41
4:Q:104:ILE:O	4:Q:106:ALA:N	2.54	0.41
4:Q:80:LEU:HA	4:Q:89:THR:O	2.21	0.41
3:C:277:LYS:HE3	7:G:31:ARG:NE	2.35	0.41
3:C:281:LYS:O	3:C:284:ALA:HB3	2.21	0.41
5:E:15:PHE:HA	5:E:15:PHE:HD2	1.65	0.41
5:E:23:ILE:O	5:E:27:LYS:HG2	2.21	0.41
6:F:26:LEU:HA	6:F:26:LEU:HD13	1.77	0.41
3:C:275:LYS:HE3	4:D:17:GLN:HB2	2.02	0.41
12:D:306:OPC:HBL1	12:D:306:OPC:CAP	2.44	0.41
8:H:24:TRP:CE3	8:H:25:GLY:N	2.89	0.41
1:N:156:GLU:HG3	1:N:166:SER:OG	2.21	0.41
2:O:106:LEU:HD21	13:O:1201:CLA:C15	2.48	0.41
7:T:11:LEU:HA	7:T:14:VAL:HG11	2.03	0.41
3:P:192:ASN:O	3:P:193:VAL:HG23	2.21	0.41
2:O:126:ARG:HA	2:O:127:PRO:HD2	1.78	0.41
1:A:142:GLN:HE21	2:B:67:ASN:N	2.16	0.41
3:C:3:PHE:HE1	3:C:117:GLY:CA	2.33	0.41
6:S:25:LEU:HD23	6:S:25:LEU:C	2.40	0.41
2:O:75:ILE:O	2:O:77:PRO:CD	2.67	0.41
6:F:25:LEU:C	6:F:25:LEU:HD23	2.41	0.41
3:C:10:PRO:HG2	3:C:11:PRO:CD	2.49	0.41
1:N:17:LEU:C	1:N:17:LEU:HD13	2.41	0.41
3:C:20:ILE:HB	3:C:240:PHE:CZ	2.56	0.41
3:C:66:SER:O	3:C:68:VAL:HG13	2.21	0.41
3:P:270:LEU:HG	3:P:274:LEU:HD11	2.03	0.41
3:C:6:GLN:HG3	3:C:106:TYR:O	2.20	0.41
3:P:4:TRP:HE3	3:P:4:TRP:N	2.18	0.41
3:C:140:LYS:HA	3:C:140:LYS:HD3	1.96	0.41
4:Q:44:PRO:HA	4:Q:45:PRO:HD2	1.51	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:C:213:ALA:C	3:C:215:PRO:HD2	2.40	0.41
3:C:131:VAL:CG2	3:C:132:LEU:N	2.83	0.41
5:E:2:ILE:O	5:E:2:ILE:HD12	2.21	0.41
1:A:14:ILE:HG23	1:A:15:GLN:N	2.31	0.41
15:R:1101:BCR:C14	6:S:16:ILE:HG21	2.51	0.41
4:D:92:VAL:O	4:D:100:ARG:HB2	2.21	0.41
4:Q:123:LYS:HE3	4:Q:125:LYS:CE	2.51	0.41
2:B:36:LEU:C	2:B:38:TYR:N	2.74	0.41
2:B:42:VAL:CG1	7:G:28:GLN:NE2	2.83	0.41
1:N:190:VAL:HG11	10:N:1304:TDS:OAK	2.21	0.41
3:C:231:LEU:HD21	3:C:233:ASN:O	2.20	0.41
2:B:152:ASP:OD2	2:B:152:ASP:C	2.58	0.41
1:A:154:VAL:N	1:A:155:PRO:HD2	2.36	0.41
3:C:52:ILE:HD12	3:C:127:ILE:HD12	2.03	0.41
1:N:23:SER:O	1:N:25:TYR:CZ	2.74	0.41
3:P:130:PRO:O	3:P:130:PRO:HG2	2.21	0.41
3:P:4:TRP:CA	3:P:4:TRP:CE3	3.03	0.41
1:N:112:LYS:CB	1:N:113:PRO:CD	2.95	0.40
1:N:28:PRO:HD2	2:O:32:TRP:CA	2.51	0.40
4:Q:15:ARG:HH22	4:Q:17:GLN:HG2	1.86	0.40
4:Q:64:VAL:HG13	4:Q:69:PHE:CE2	2.56	0.40
4:Q:84:LEU:O	4:Q:87:ASP:OD1	2.38	0.40
2:B:118:ASN:HB3	2:B:121:GLN:HG3	2.02	0.40
8:H:22:VAL:HG12	8:H:26:ARG:HE	1.85	0.40
3:C:188:ASP:O	3:C:189:GLU:CB	2.68	0.40
4:D:160:ILE:HG22	4:D:162:LEU:HD13	2.04	0.40
2:B:84:VAL:CG1	2:B:101:MET:HB2	2.50	0.40
3:C:127:ILE:HG22	3:C:129:PHE:CE1	2.57	0.40
3:C:161:TYR:C	3:C:163:THR:N	2.74	0.40
1:N:25:TYR:HB2	1:N:26:VAL:H	1.51	0.40
1:N:50:THR:O	1:N:54:MET:HG3	2.20	0.40
1:N:102:PHE:HB3	5:R:18:ILE:CD1	2.49	0.40
3:P:275:LYS:CE	4:Q:20:ASN:HD22	2.34	0.40
3:C:275:LYS:HG3	4:D:20:ASN:HB3	2.02	0.40
3:C:192:ASN:O	3:C:193:VAL:HG23	2.21	0.40
10:A:304:TDS:HBC1	2:B:85:PHE:CB	2.51	0.40
2:B:149:LEU:HD23	4:Q:129:HIS:CD2	2.56	0.40
3:C:161:TYR:C	3:C:163:THR:H	2.25	0.40
4:D:59:LYS:HD3	4:D:79:VAL:HB	2.03	0.40
6:S:3:GLU:HA	6:S:6:LEU:HD12	2.02	0.40
3:C:79:PRO:O	3:C:80:GLU:C	2.60	0.40
1:A:34:TYR:CD2	1:A:34:TYR:N	2.87	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:D:95:SER:C	4:D:97:GLU:N	2.75	0.40
4:Q:131:SER:CB	4:Q:144:ALA:HB2	2.51	0.40
4:Q:66:VAL:CG2	4:Q:160:ILE:HG13	2.46	0.40
4:Q:165:TRP:CE2	4:Q:178:TRP:CH2	3.10	0.40
4:Q:138:ARG:CD	4:Q:171:ARG:HD3	2.51	0.40
4:Q:83:GLY:CA	4:Q:87:ASP:O	2.69	0.40
2:B:126:ARG:HG2	2:B:128:VAL:CG2	2.51	0.40
2:B:57:LEU:HD13	7:G:15:PHE:CD2	2.56	0.40
7:G:9:LEU:HG	7:G:10:VAL:N	2.36	0.40
2:B:22:MET:HE2	2:B:22:MET:HA	2.03	0.40
3:C:185:LYS:CD	3:C:217:LEU:HD11	2.51	0.40
3:C:93:GLU:HG3	3:C:94:LEU:H	1.86	0.40
4:D:50:VAL:HG12	4:D:164:PRO:CB	2.49	0.40
1:A:134:THR:HG23	10:A:304:TDS:HAQ2	2.03	0.40
4:Q:111:LEU:HD23	4:Q:129:HIS:CE1	2.57	0.40
3:C:226:LYS:HB3	3:C:227:ALA:H	1.69	0.40
1:N:83:ARG:NH2	2:O:61:MET:O	2.54	0.40
1:A:116:LEU:CD1	1:A:116:LEU:H	2.18	0.40
1:N:189:PHE:CD1	1:N:193:TRP:CZ3	3.09	0.40
2:O:135:PHE:O	2:O:139:VAL:HG23	2.22	0.40
3:P:59:GLN:OE1	3:P:67:LYS:HE3	2.21	0.40
3:C:14:ARG:NH2	3:C:150:HIS:CD2	2.89	0.40
5:R:5:ALA:HB3	6:S:5:MET:HB2	2.01	0.40
8:U:17:TRP:O	8:U:21:MET:HG3	2.22	0.40
3:C:61:VAL:O	3:C:61:VAL:HG23	2.21	0.40
4:Q:96:LYS:O	4:Q:96:LYS:HD2	2.22	0.40
3:P:76:LEU:HG	3:P:78:LEU:HD21	2.03	0.40
1:N:163:VAL:HG12	1:N:164:LEU:N	2.36	0.40
4:D:102:TYR:OH	4:D:135:GLU:HG3	2.21	0.40
4:D:149:ALA:HB1	4:D:178:TRP:CD1	2.57	0.40
4:Q:155:VAL:HA	4:Q:161:VAL:N	2.24	0.40
7:G:21:LEU:HD13	7:G:21:LEU:C	2.41	0.40
7:G:25:ALA:HA	7:G:28:GLN:OE1	2.22	0.40
3:P:188:ASP:O	3:P:189:GLU:CB	2.69	0.40
3:P:193:VAL:HG12	3:P:195:TYR:CE1	2.57	0.40
3:C:109:GLY:O	3:C:111:ASP:N	2.55	0.40
6:S:32:ALA:O	6:S:34:LYS:N	2.55	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	200/215 (93%)	152 (76%)	29 (14%)	19 (10%)	1	4
1	N	200/215 (93%)	155 (78%)	32 (16%)	13 (6%)	2	11
2	B	136/160 (85%)	80 (59%)	31 (23%)	25 (18%)	0	1
2	O	136/160 (85%)	67 (49%)	38 (28%)	31 (23%)	0	0
3	C	284/289 (98%)	206 (72%)	47 (16%)	31 (11%)	1	3
3	P	284/289 (98%)	213 (75%)	44 (16%)	27 (10%)	1	4
4	D	166/179 (93%)	123 (74%)	25 (15%)	18 (11%)	1	3
4	Q	166/179 (93%)	85 (51%)	49 (30%)	32 (19%)	0	0
5	E	30/32 (94%)	22 (73%)	6 (20%)	2 (7%)	2	10
5	R	30/32 (94%)	23 (77%)	5 (17%)	2 (7%)	2	10
6	F	31/35 (89%)	25 (81%)	5 (16%)	1 (3%)	6	33
6	S	33/35 (94%)	25 (76%)	5 (15%)	3 (9%)	1	5
7	G	21/37 (57%)	15 (71%)	5 (24%)	1 (5%)	4	20
7	T	25/37 (68%)	18 (72%)	6 (24%)	1 (4%)	5	25
8	H	25/29 (86%)	18 (72%)	7 (28%)	0	100	100
8	U	25/29 (86%)	19 (76%)	6 (24%)	0	100	100
All	All	1792/1952 (92%)	1246 (70%)	340 (19%)	206 (12%)	1	3

All (206) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	25	TYR
1	A	28	PRO
1	A	111	LYS
1	A	160	VAL
1	A	173	SER
1	A	211	ILE
2	B	22	MET
2	B	25	ASN

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Mol	Chain	Res	Type
2	B	32	TRP
2	B	33	PRO
2	B	34	ASN
2	B	35	ASP
2	B	72	PRO
2	B	75	ILE
2	B	76	LEU
2	B	149	LEU
2	B	154	THR
3	C	54	TYR
3	C	92	GLU
3	C	99	GLY
3	C	110	GLN
3	C	189	GLU
3	C	196	GLN
3	C	203	SER
3	C	227	ALA
3	C	232	THR
3	C	276	LYS
4	D	66	VAL
4	D	85	LYS
4	D	110	HIS
1	N	19	ASP
1	N	21	VAL
1	N	22	THR
1	N	30	VAL
1	N	32	ILE
1	N	111	LYS
2	O	25	ASN
2	O	26	TYR
2	O	30	PRO
2	O	32	TRP
2	O	33	PRO
2	O	34	ASN
2	O	35	ASP
2	O	65	PRO
2	O	72	PRO
2	O	73	LEU
2	O	115	GLU
2	O	121	GLN
2	O	123	PRO
2	O	124	PHE

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Mol	Chain	Res	Type
2	O	150	PRO
2	O	151	LEU
3	P	54	TYR
3	P	92	GLU
3	P	99	GLY
3	P	110	GLN
3	P	189	GLU
3	P	196	GLN
3	P	203	SER
3	P	227	ALA
3	P	232	THR
3	P	276	LYS
4	Q	46	SER
4	Q	57	LYS
4	Q	73	HIS
4	Q	74	ASN
4	Q	92	VAL
4	Q	106	ALA
4	Q	124	PHE
4	Q	145	PRO
4	Q	153	ALA
4	Q	164	PRO
1	A	14	ILE
1	A	20	ASP
1	A	27	PRO
1	A	75	GLU
1	A	106	LEU
2	B	19	ALA
2	B	71	THR
2	B	78	GLU
2	B	93	ASN
2	B	115	GLU
2	B	128	VAL
2	B	145	ILE
2	B	146	GLY
2	B	153	LYS
3	C	56	THR
3	C	94	LEU
3	C	188	ASP
3	C	202	ASP
3	C	230	ALA
4	D	14	GLY

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Mol	Chain	Res	Type
4	D	20	ASN
4	D	75	ALA
4	D	106	ALA
4	D	109	THR
7	G	29	TYR
1	N	16	ALA
1	N	110	PHE
1	N	177	GLN
1	N	190	VAL
2	O	20	LYS
2	O	93	ASN
2	O	128	VAL
2	O	147	ALA
2	O	148	ALA
2	O	149	LEU
2	O	153	LYS
3	P	56	THR
3	P	188	ASP
3	P	202	ASP
3	P	230	ALA
4	Q	14	GLY
4	Q	56	ALA
4	Q	81	VAL
4	Q	91	ILE
4	Q	93	VAL
4	Q	104	ILE
4	Q	105	ASN
4	Q	147	SER
4	Q	157	ASP
5	R	25	ALA
6	S	33	GLU
7	T	29	TYR
1	A	112	LYS
1	A	140	TRP
1	A	190	VAL
2	B	92	PRO
3	C	23	ALA
3	C	127	ILE
3	C	220	SER
3	C	233	ASN
4	D	68	LYS
4	D	99	ILE

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Mol	Chain	Res	Type
4	D	141	ARG
5	E	25	ALA
1	N	205	MET
2	O	63	GLY
2	O	64	GLU
2	O	92	PRO
2	O	114	ILE
2	O	118	ASN
3	P	94	LEU
3	P	193	VAL
3	P	220	SER
3	P	233	ASN
4	Q	48	GLY
4	Q	175	LYS
1	A	15	GLN
1	A	205	MET
2	B	30	PRO
2	B	79	TRP
3	C	87	GLU
3	C	130	PRO
3	C	187	GLU
3	C	193	VAL
4	D	56	ALA
2	O	75	ILE
3	P	23	ALA
3	P	87	GLU
3	P	91	PRO
3	P	127	ILE
3	P	130	PRO
3	P	187	GLU
4	Q	13	MET
4	Q	20	ASN
4	Q	45	PRO
4	Q	67	SER
4	Q	143	PRO
1	A	24	LYS
1	A	105	TYR
2	B	114	ILE
3	C	221	GLU
4	D	47	GLY
4	D	50	VAL
4	D	101	ASP

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Mol	Chain	Res	Type
4	D	143	PRO
2	O	76	LEU
3	P	224	ALA
4	Q	50	VAL
4	Q	98	ALA
4	Q	151	CYS
4	Q	160	ILE
6	S	30	GLN
1	A	21	VAL
3	C	58	LEU
3	C	191	GLY
3	C	224	ALA
3	C	282	VAL
4	D	13	MET
1	N	26	VAL
4	Q	102	TYR
3	C	31	PRO
5	E	20	VAL
3	P	282	VAL
3	P	191	GLY
6	S	29	ILE
3	C	91	PRO
6	F	29	ILE
2	O	109	ILE
5	R	20	VAL
2	B	123	PRO
3	C	117	GLY
4	D	145	PRO
4	Q	64	VAL
1	N	211	ILE

### 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	172/184 (94%)	147 (86%)	25 (14%)	5	22
1	N	172/184 (94%)	153 (89%)	19 (11%)	9	34

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	117/136 (86%)	93 (80%)	24 (20%)	2	9
2	O	117/136 (86%)	98 (84%)	19 (16%)	3	17
3	C	240/243 (99%)	207 (86%)	33 (14%)	5	24
3	P	240/243 (99%)	206 (86%)	34 (14%)	5	22
4	D	136/146 (93%)	113 (83%)	23 (17%)	3	15
4	Q	136/146 (93%)	113 (83%)	23 (17%)	3	15
5	E	25/25 (100%)	15 (60%)	10 (40%)	0	0
5	R	25/25 (100%)	14 (56%)	11 (44%)	0	0
6	F	25/27 (93%)	22 (88%)	3 (12%)	7	30
6	S	27/27 (100%)	23 (85%)	4 (15%)	4	21
7	G	17/28 (61%)	9 (53%)	8 (47%)	0	0
7	T	21/28 (75%)	12 (57%)	9 (43%)	0	0
8	H	22/24 (92%)	18 (82%)	4 (18%)	2	13
8	U	22/24 (92%)	18 (82%)	4 (18%)	2	13
All	All	1514/1626 (93%)	1261 (83%)	253 (17%)	3	16

All (253) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ILE
1	A	15	GLN
1	A	19	ASP
1	A	20	ASP
1	A	24	LYS
1	A	26	VAL
1	A	28	PRO
1	A	36	LEU
1	A	47	GLN
1	A	63	THR
1	A	74	ASN
1	A	75	GLU
1	A	95	LEU
1	A	103	ARG
1	A	115	GLU
1	A	116	LEU
1	A	140	TRP
1	A	150	ILE

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Mol	Chain	Res	Type
1	A	164	LEU
1	A	167	ASP
1	A	173	SER
1	A	177	GLN
1	A	193	TRP
1	A	200	LEU
1	A	207	ARG
2	B	20	LYS
2	B	22	MET
2	B	25	ASN
2	B	26	TYR
2	B	27	TYR
2	B	29	GLU
2	B	32	TRP
2	B	33	PRO
2	B	35	ASP
2	B	36	LEU
2	B	38	TYR
2	B	62	VAL
2	B	74	GLU
2	B	78	GLU
2	B	79	TRP
2	B	80	TYR
2	B	81	LEU
2	B	90	SER
2	B	95	LEU
2	B	96	LEU
2	B	116	ASN
2	B	121	GLN
2	B	124	PHE
2	B	149	LEU
3	C	3	PHE
3	C	4	TRP
3	C	9	TYR
3	C	14	ARG
3	C	17	THR
3	C	19	ARG
3	C	27	LEU
3	C	51	LYS
3	C	83	LYS
3	C	93	GLU
3	C	94	LEU

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Mol	Chain	Res	Type
3	C	98	VAL
3	C	100	ASP
3	C	104	GLN
3	C	119	LEU
3	C	122	GLU
3	C	123	GLN
3	C	131	VAL
3	C	155	ARG
3	C	165	GLU
3	C	182	LYS
3	C	205	LYS
3	C	211	ILE
3	C	221	GLU
3	C	231	LEU
3	C	232	THR
3	C	233	ASN
3	C	249	LEU
3	C	258	MET
3	C	264	LEU
3	C	271	MET
3	C	276	LYS
3	C	281	LYS
4	D	13	MET
4	D	15	ARG
4	D	16	ARG
4	D	22	LEU
4	D	62	ASN
4	D	63	ASN
4	D	82	GLN
4	D	92	VAL
4	D	93	VAL
4	D	100	ARG
4	D	101	ASP
4	D	102	TYR
4	D	126	CYS
4	D	138	ARG
4	D	139	VAL
4	D	154	THR
4	D	158	ASP
4	D	161	VAL
4	D	169	ASP
4	D	170	PHE

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Mol	Chain	Res	Type
4	D	174	GLU
4	D	175	LYS
4	D	177	TRP
5	E	1	MET
5	E	2	ILE
5	E	8	TYR
5	E	11	PHE
5	E	12	ILE
5	E	14	LEU
5	E	15	PHE
5	E	16	PHE
5	E	29	ILE
5	E	30	LYS
6	F	7	TYR
6	F	26	LEU
6	F	30	GLN
7	G	9	LEU
7	G	11	LEU
7	G	13	LEU
7	G	14	VAL
7	G	18	LEU
7	G	22	PHE
7	G	26	TYR
7	G	31	ARG
8	H	8	TRP
8	H	16	THR
8	H	21	MET
8	H	24	TRP
1	N	13	GLU
1	N	17	LEU
1	N	25	TYR
1	N	29	HIS
1	N	31	ASN
1	N	47	GLN
1	N	63	THR
1	N	95	LEU
1	N	103	ARG
1	N	110	PHE
1	N	112	LYS
1	N	116	LEU
1	N	140	TRP
1	N	150	ILE

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Mol	Chain	Res	Type
1	N	163	VAL
1	N	164	LEU
1	N	167	ASP
1	N	193	TRP
1	N	200	LEU
2	O	24	HIS
2	O	26	TYR
2	O	27	TYR
2	O	30	PRO
2	O	32	TRP
2	O	33	PRO
2	O	35	ASP
2	O	36	LEU
2	O	57	LEU
2	O	64	GLU
2	O	65	PRO
2	O	74	GLU
2	O	90	SER
2	O	92	PRO
2	O	95	LEU
2	O	96	LEU
2	O	108	LEU
2	O	116	ASN
2	O	152	ASP
3	P	3	PHE
3	P	4	TRP
3	P	14	ARG
3	P	17	THR
3	P	19	ARG
3	P	27	LEU
3	P	51	LYS
3	P	83	LYS
3	P	93	GLU
3	P	94	LEU
3	P	98	VAL
3	P	100	ASP
3	P	104	GLN
3	P	119	LEU
3	P	122	GLU
3	P	123	GLN
3	P	129	PHE
3	P	131	VAL

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Mol	Chain	Res	Type
3	P	155	ARG
3	P	165	GLU
3	P	182	LYS
3	P	205	LYS
3	P	206	THR
3	P	211	ILE
3	P	221	GLU
3	P	231	LEU
3	P	232	THR
3	P	233	ASN
3	P	249	LEU
3	P	258	MET
3	P	264	LEU
3	P	271	MET
3	P	276	LYS
3	P	281	LYS
4	Q	13	MET
4	Q	15	ARG
4	Q	16	ARG
4	Q	22	LEU
4	Q	58	ASP
4	Q	59	LYS
4	Q	63	ASN
4	Q	64	VAL
4	Q	71	GLU
4	Q	77	ASP
4	Q	94	GLU
4	Q	97	GLU
4	Q	132	GLN
4	Q	134	ASP
4	Q	138	ARG
4	Q	156	GLN
4	Q	158	ASP
4	Q	159	ASN
4	Q	164	PRO
4	Q	170	PHE
4	Q	174	GLU
4	Q	175	LYS
4	Q	178	TRP
5	R	1	MET
5	R	2	ILE
5	R	8	TYR

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Mol	Chain	Res	Type
5	R	11	PHE
5	R	12	ILE
5	R	14	LEU
5	R	15	PHE
5	R	16	PHE
5	R	27	LYS
5	R	29	ILE
5	R	30	LYS
6	S	7	TYR
6	S	26	LEU
6	S	30	GLN
6	S	34	LYS
7	T	9	LEU
7	T	11	LEU
7	T	13	LEU
7	T	14	VAL
7	T	18	LEU
7	T	22	PHE
7	T	26	TYR
7	T	31	ARG
7	T	33	ASN
8	U	8	TRP
8	U	16	THR
8	U	21	MET
8	U	24	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	29	HIS
1	A	47	GLN
1	A	142	GLN
2	B	25	ASN
2	B	86	GLN
2	B	116	ASN
2	B	118	ASN
2	B	121	GLN
3	C	7	GLN
3	C	60	GLN
3	C	125	GLN
3	C	150	HIS

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Mol	Chain	Res	Type
3	C	159	GLN
3	C	223	GLN
3	C	236	ASN
3	C	250	GLN
3	C	269	GLN
3	C	278	GLN
4	D	82	GLN
4	D	132	GLN
4	D	152	HIS
7	G	28	GLN
8	H	27	ASN
1	N	47	GLN
1	N	142	GLN
1	N	209	GLN
2	O	86	GLN
2	O	116	ASN
2	O	121	GLN
2	O	122	ASN
3	P	7	GLN
3	P	60	GLN
3	P	125	GLN
3	P	159	GLN
3	P	223	GLN
3	P	234	ASN
3	P	236	ASN
3	P	250	GLN
4	Q	20	ASN
4	Q	63	ASN
4	Q	82	GLN
4	Q	105	ASN
4	Q	122	ASN
4	Q	132	GLN
7	T	27	GLN
8	U	27	ASN

### 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 ⓘ

22 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	$\# Z  > 2$
9	HEM	A	301	1	49,50,50	3.94	15 (30%)	46,82,82	1.60	10 (21%)
9	HEM	A	302	1	49,50,50	3.38	16 (32%)	46,82,82	2.86	11 (23%)
9	HEM	A	303	1,16	49,50,50	4.10	16 (32%)	46,82,82	2.21	14 (30%)
10	TDS	A	304	-	31,31,31	2.60	6 (19%)	38,40,40	4.54	12 (31%)
11	PL9	A	305	-	55,55,55	3.53	21 (38%)	69,69,69	2.45	21 (30%)
13	CLA	B	201	-	73,73,73	1.74	14 (19%)	95,113,113	2.18	23 (24%)
12	OPC	B	307	-	51,53,54	1.37	6 (11%)	55,61,64	1.06	6 (10%)
9	HEM	C	301	3	49,50,50	2.93	17 (34%)	46,82,82	1.49	9 (19%)
14	FES	D	200	4	0,4,4	0.00	-	0,4,4	0.00	-
12	OPC	D	306	-	51,53,54	1.38	6 (11%)	55,61,64	1.14	4 (7%)
15	BCR	E	101	-	41,41,41	2.09	8 (19%)	56,56,56	2.44	23 (41%)
10	TDS	N	1304	-	31,31,31	2.58	6 (19%)	38,40,40	4.27	13 (34%)
12	OPC	N	1306	-	51,53,54	1.37	6 (11%)	55,61,64	1.27	8 (14%)
9	HEM	N	301	1	49,50,50	3.32	17 (34%)	46,82,82	1.67	9 (19%)
9	HEM	N	302	1	49,50,50	2.95	17 (34%)	46,82,82	2.88	13 (28%)
9	HEM	N	303	1,16	49,50,50	4.25	19 (38%)	46,82,82	1.82	8 (17%)
13	CLA	O	1201	-	73,73,73	1.76	13 (17%)	95,113,113	2.23	22 (23%)
9	HEM	P	301	3	49,50,50	3.36	19 (38%)	46,82,82	1.49	7 (15%)
14	FES	Q	1200	4	0,4,4	0.00	-	0,4,4	0.00	-
11	PL9	Q	1305	-	55,55,55	3.52	21 (38%)	69,69,69	2.56	22 (31%)
12	OPC	Q	1307	-	51,53,54	1.37	6 (11%)	55,61,64	1.11	5 (9%)
15	BCR	R	1101	-	41,41,41	2.03	8 (19%)	56,56,56	2.45	22 (39%)

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
9	HEM	A	301	1	-	0/14/114/114	0/0/8/8
9	HEM	A	302	1	-	0/14/114/114	0/0/8/8
9	HEM	A	303	1,16	-	0/14/114/114	0/0/8/8
10	TDS	A	304	-	-	0/16/17/17	0/0/2/2
11	PL9	A	305	-	-	1/53/73/73	0/1/1/1
13	CLA	B	201	-	2/2/20/25	0/37/135/135	0/0/9/9
12	OPC	B	307	-	-	2/55/57/60	0/0/0/0
9	HEM	C	301	3	-	0/14/114/114	0/0/8/8
14	FES	D	200	4	-	0/0/4/4	0/0/1/1
12	OPC	D	306	-	-	3/55/57/60	0/0/0/0
15	BCR	E	101	-	-	0/29/63/63	0/2/2/2
10	TDS	N	1304	-	-	0/16/17/17	0/0/2/2
12	OPC	N	1306	-	-	3/55/57/60	0/0/0/0
9	HEM	N	301	1	-	0/14/114/114	0/0/8/8
9	HEM	N	302	1	-	0/14/114/114	0/0/8/8
9	HEM	N	303	1,16	-	0/14/114/114	0/0/8/8
13	CLA	O	1201	-	3/3/20/25	0/37/135/135	0/0/9/9
9	HEM	P	301	3	-	0/14/114/114	0/0/8/8
14	FES	Q	1200	4	-	0/0/4/4	0/0/1/1
11	PL9	Q	1305	-	-	1/53/73/73	0/1/1/1
12	OPC	Q	1307	-	-	2/55/57/60	0/0/0/0
15	BCR	R	1101	-	-	0/29/63/63	0/2/2/2

All (257) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	301	HEM	C2D-C1D	22.13	1.50	1.44
9	A	302	HEM	C2D-C1D	18.41	1.49	1.44
9	N	303	HEM	C2D-C1D	17.62	1.49	1.44
9	N	301	HEM	C2D-C1D	17.05	1.48	1.44
9	A	303	HEM	C2D-C1D	15.67	1.48	1.44
9	N	303	HEM	C3D-C4D	15.58	1.48	1.44
9	A	303	HEM	C3D-C4D	15.36	1.48	1.44
9	P	301	HEM	C2D-C1D	14.64	1.48	1.44
9	N	302	HEM	C2D-C1D	13.08	1.47	1.44
9	C	301	HEM	C2D-C1D	10.77	1.47	1.44
11	A	305	PL9	C13-C14	10.28	1.53	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	A	305	PL9	C33-C34	10.21	1.53	1.32
11	A	305	PL9	C18-C19	10.20	1.53	1.32
11	Q	1305	PL9	C13-C14	10.19	1.53	1.32
11	Q	1305	PL9	C33-C34	10.16	1.53	1.32
11	Q	1305	PL9	C18-C19	10.15	1.53	1.32
9	A	301	HEM	C4A-C3A	9.29	1.51	1.40
10	N	1304	TDS	CAL-CAM	9.15	1.53	1.40
10	A	304	TDS	CAL-CAM	9.14	1.53	1.40
9	C	301	HEM	C3D-C4D	8.74	1.46	1.44
9	P	301	HEM	C3D-C4D	8.19	1.46	1.44
9	A	303	HEM	C2B-C1B	-7.78	1.42	1.44
10	A	304	TDS	CAQ-CAP	-7.15	1.39	1.51
10	N	1304	TDS	CAQ-CAP	-7.10	1.39	1.51
9	N	303	HEM	C2B-C1B	-7.00	1.42	1.44
9	P	301	HEM	C4A-C3A	6.99	1.48	1.40
9	N	301	HEM	C4A-C3A	6.93	1.48	1.40
11	A	305	PL9	C7-C8	-6.92	1.39	1.50
11	Q	1305	PL9	C7-C8	-6.90	1.39	1.50
9	N	302	HEM	C2B-C1B	-6.84	1.42	1.44
9	N	303	HEM	C4A-C3A	6.79	1.48	1.40
9	A	303	HEM	C4A-C3A	6.69	1.48	1.40
11	Q	1305	PL9	C48-C49	6.54	1.53	1.32
15	E	101	BCR	C26-C25	6.52	1.44	1.34
11	A	305	PL9	C48-C49	6.51	1.53	1.32
9	A	302	HEM	C4A-C3A	6.34	1.48	1.40
9	C	301	HEM	C4A-C3A	6.31	1.48	1.40
15	E	101	BCR	C30-C25	6.23	1.62	1.53
15	R	1101	BCR	C26-C25	6.22	1.44	1.34
13	B	201	CLA	C3A-C2A	-5.85	1.36	1.54
9	A	303	HEM	C3D-C2D	-5.75	1.33	1.43
11	Q	1305	PL9	C3-C4	-5.53	1.39	1.49
11	A	305	PL9	C3-C4	-5.44	1.39	1.49
9	N	303	HEM	C3D-C2D	-5.43	1.34	1.43
13	O	1201	CLA	C3B-C4B	5.41	1.48	1.40
13	B	201	CLA	C4D-C3D	-5.38	1.35	1.41
13	B	201	CLA	C3B-C4B	5.31	1.48	1.40
15	R	1101	BCR	C30-C25	5.13	1.61	1.53
10	A	304	TDS	CAM-CAN	5.11	1.53	1.42
10	N	1304	TDS	CAM-CAN	5.10	1.53	1.42
9	C	301	HEM	C3C-CAC	5.10	1.56	1.40
9	A	303	HEM	C3B-C2B	-5.10	1.34	1.43
13	O	1201	CLA	C3A-C2A	-5.06	1.39	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	N	301	HEM	C3B-C2B	-5.05	1.34	1.43
9	P	301	HEM	C3C-CAC	5.03	1.56	1.40
9	A	303	HEM	CBB-CAB	4.91	1.57	1.28
9	N	302	HEM	C3C-CAC	4.91	1.55	1.40
11	A	305	PL9	C11-C9	-4.90	1.39	1.51
11	A	305	PL9	C41-C39	-4.90	1.39	1.51
11	Q	1305	PL9	C11-C9	-4.89	1.39	1.51
9	N	303	HEM	CBB-CAB	4.87	1.57	1.28
11	Q	1305	PL9	C41-C39	-4.87	1.39	1.51
11	Q	1305	PL9	C6-C1	-4.84	1.39	1.48
13	O	1201	CLA	C4D-C3D	-4.69	1.35	1.41
11	A	305	PL9	C6-C1	-4.69	1.40	1.48
9	N	303	HEM	C3B-C2B	-4.64	1.35	1.43
15	R	1101	BCR	C1-C6	4.60	1.60	1.53
9	N	302	HEM	C4A-C3A	4.58	1.45	1.40
9	N	303	HEM	C3B-CAB	4.58	1.54	1.40
9	A	303	HEM	C3B-CAB	4.57	1.54	1.40
9	A	302	HEM	C3B-C4B	4.56	1.49	1.44
9	A	303	HEM	C3C-CAC	4.55	1.54	1.40
9	P	301	HEM	CBC-CAC	4.53	1.55	1.28
9	A	303	HEM	CBC-CAC	4.53	1.55	1.28
9	C	301	HEM	CBC-CAC	4.52	1.55	1.28
15	E	101	BCR	C1-C6	4.50	1.60	1.53
9	A	303	HEM	C3B-C4B	4.48	1.49	1.44
9	P	301	HEM	CHA-C4D	4.39	1.42	1.35
9	C	301	HEM	CHA-C4D	4.35	1.42	1.35
9	N	302	HEM	CBC-CAC	4.34	1.54	1.28
9	A	301	HEM	C2B-C1B	-4.34	1.43	1.44
9	P	301	HEM	C3B-C2B	-4.25	1.36	1.43
9	N	303	HEM	C3C-CAC	4.22	1.53	1.40
9	A	302	HEM	C3C-CAC	4.22	1.53	1.40
9	N	303	HEM	CBC-CAC	4.21	1.53	1.28
9	N	301	HEM	C3C-CAC	4.21	1.53	1.40
9	N	303	HEM	C3C-C2C	-4.19	1.36	1.43
9	A	302	HEM	CBC-CAC	4.19	1.53	1.28
13	O	1201	CLA	O2D-CGD	4.09	1.44	1.33
9	N	303	HEM	C3B-C4B	4.09	1.49	1.44
10	A	304	TDS	OAO-CAP	-4.07	1.36	1.38
9	A	301	HEM	CBC-CAC	4.04	1.52	1.28
13	B	201	CLA	O2D-CGD	4.03	1.43	1.33
13	O	1201	CLA	C2-C3	4.00	1.41	1.32
12	Q	1307	OPC	CAV-CAW	3.99	1.53	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	P	301	HEM	C2B-C1B	3.97	1.45	1.44
9	N	302	HEM	C3B-C2B	-3.96	1.36	1.43
9	A	301	HEM	C3C-CAC	3.96	1.52	1.40
12	D	306	OPC	CAV-CAW	3.95	1.53	1.31
12	N	1306	OPC	CAV-CAW	3.94	1.53	1.31
10	A	304	TDS	OAK-CAL	3.93	1.43	1.37
9	N	301	HEM	CBC-CAC	3.92	1.51	1.28
12	B	307	OPC	CAV-CAW	3.92	1.53	1.31
11	Q	1305	PL9	C41-C42	-3.92	1.39	1.53
9	A	301	HEM	C3B-C2B	-3.91	1.36	1.43
11	A	305	PL9	C41-C42	-3.91	1.39	1.53
10	N	1304	TDS	OAO-CAP	-3.87	1.36	1.38
10	N	1304	TDS	OAK-CAL	3.87	1.43	1.37
9	A	303	HEM	C3C-C2C	-3.83	1.37	1.43
11	A	305	PL9	C47-C48	-3.79	1.39	1.50
11	A	305	PL9	C27-C28	-3.79	1.39	1.50
11	Q	1305	PL9	C47-C48	-3.78	1.39	1.50
11	A	305	PL9	C22-C23	-3.78	1.39	1.50
11	Q	1305	PL9	C27-C28	-3.77	1.39	1.50
11	Q	1305	PL9	C22-C23	-3.77	1.39	1.50
11	A	305	PL9	C42-C43	-3.71	1.39	1.50
9	P	301	HEM	C3B-C4B	3.70	1.48	1.44
15	R	1101	BCR	C38-C26	3.69	1.57	1.51
11	Q	1305	PL9	C42-C43	-3.68	1.39	1.50
9	N	302	HEM	FE-NB	3.62	2.11	1.97
9	N	302	HEM	C3B-C4B	3.62	1.48	1.44
13	O	1201	CLA	C1B-C2B	3.59	1.44	1.40
9	C	301	HEM	C3B-C2B	-3.55	1.37	1.43
9	N	301	HEM	CHC-C1C	3.54	1.42	1.36
9	N	301	HEM	C4D-ND	-3.53	1.32	1.39
9	N	302	HEM	C3D-C4D	3.51	1.45	1.44
12	D	306	OPC	CBP-CBQ	-3.51	1.39	1.52
12	Q	1307	OPC	CBP-CBQ	-3.50	1.39	1.52
12	B	307	OPC	CBP-CBQ	-3.50	1.39	1.52
11	Q	1305	PL9	C51-C49	-3.49	1.39	1.50
9	A	302	HEM	CHC-C1C	3.49	1.42	1.36
11	A	305	PL9	C51-C49	-3.45	1.39	1.50
12	N	1306	OPC	CBP-CBQ	-3.45	1.39	1.52
9	A	302	HEM	FE-NB	3.44	2.10	1.97
9	A	302	HEM	C3B-C2B	-3.41	1.37	1.43
11	A	305	PL9	C43-C44	3.40	1.39	1.32
15	E	101	BCR	C38-C26	3.39	1.56	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	N	301	HEM	C3C-C2C	-3.39	1.37	1.43
9	P	301	HEM	C3C-C2C	-3.38	1.37	1.43
13	O	1201	CLA	C1A-NA	3.37	1.39	1.32
13	B	201	CLA	C2-C3	3.36	1.39	1.32
11	Q	1305	PL9	C43-C44	3.35	1.39	1.32
9	A	301	HEM	FE-NA	3.35	2.06	1.92
11	Q	1305	PL9	C38-C39	3.34	1.39	1.32
11	A	305	PL9	C23-C24	3.34	1.39	1.32
11	Q	1305	PL9	C23-C24	3.33	1.39	1.32
11	A	305	PL9	C38-C39	3.32	1.39	1.32
9	P	301	HEM	FE-NB	3.32	2.09	1.97
11	A	305	PL9	C28-C29	3.32	1.39	1.32
9	A	301	HEM	C4D-ND	-3.30	1.32	1.39
13	B	201	CLA	O1D-CGD	3.30	1.29	1.21
15	E	101	BCR	C29-C30	3.30	1.62	1.54
9	C	301	HEM	FE-NC	3.30	2.10	1.97
9	N	301	HEM	FE-NA	3.29	2.06	1.92
11	A	305	PL9	C8-C9	3.28	1.39	1.32
12	B	307	OPC	CAQ-CAP	-3.28	1.39	1.52
12	D	306	OPC	CAQ-CAP	-3.28	1.39	1.52
9	A	302	HEM	C3C-C2C	-3.27	1.38	1.43
11	Q	1305	PL9	C8-C9	3.27	1.39	1.32
11	Q	1305	PL9	C28-C29	3.26	1.39	1.32
9	C	301	HEM	C3C-C2C	-3.25	1.38	1.43
12	N	1306	OPC	CAQ-CAP	-3.25	1.39	1.52
15	R	1101	BCR	C29-C30	3.24	1.62	1.54
12	Q	1307	OPC	CAQ-CAP	-3.24	1.39	1.52
9	P	301	HEM	C1A-C2A	3.21	1.49	1.43
9	A	301	HEM	C3C-C2C	-3.21	1.38	1.43
13	O	1201	CLA	O1D-CGD	3.21	1.29	1.21
12	D	306	OPC	CAG-CAH	-3.19	1.39	1.51
12	Q	1307	OPC	CAG-CAH	-3.17	1.39	1.51
15	R	1101	BCR	C5-C6	3.15	1.39	1.34
9	N	303	HEM	CMD-C2D	3.15	1.57	1.47
15	R	1101	BCR	C2-C1	3.15	1.62	1.54
12	B	307	OPC	CAG-CAH	-3.14	1.39	1.51
12	N	1306	OPC	CAG-CAH	-3.11	1.39	1.51
9	N	301	HEM	C3D-C2D	-3.10	1.38	1.43
9	N	301	HEM	C3B-C4B	3.08	1.48	1.44
9	A	303	HEM	CHD-C4C	3.00	1.41	1.36
9	P	301	HEM	FE-NA	2.99	2.05	1.92
9	N	302	HEM	C4D-ND	-2.99	1.33	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	N	301	HEM	FE-NB	2.98	2.08	1.97
15	E	101	BCR	C2-C1	2.94	1.61	1.54
9	A	301	HEM	C3D-C2D	-2.90	1.38	1.43
9	P	301	HEM	FE-NC	2.90	2.08	1.97
9	N	302	HEM	CHC-C1C	2.89	1.41	1.36
9	C	301	HEM	FE-NB	2.84	2.08	1.97
9	A	302	HEM	C4D-ND	-2.83	1.33	1.39
13	B	201	CLA	C1A-NA	2.83	1.38	1.32
9	A	301	HEM	CHD-C4C	2.81	1.41	1.36
11	A	305	PL9	C7-C3	2.79	1.53	1.51
9	N	301	HEM	CHD-C4C	2.79	1.41	1.36
9	N	302	HEM	C3D-C2D	-2.78	1.38	1.43
13	O	1201	CLA	CBD-CHA	2.77	1.63	1.53
15	E	101	BCR	C5-C6	2.72	1.38	1.34
9	C	301	HEM	FE-NA	2.68	2.03	1.92
9	N	302	HEM	FE-NC	2.68	2.07	1.97
9	C	301	HEM	C3B-C4B	2.65	1.47	1.44
9	N	301	HEM	C2B-C1B	2.62	1.45	1.44
9	A	302	HEM	C3B-CAB	2.60	1.48	1.40
9	C	301	HEM	C3B-CAB	2.60	1.48	1.40
9	P	301	HEM	CHC-C1C	2.56	1.41	1.36
9	A	302	HEM	C2B-C1B	-2.56	1.43	1.44
9	N	302	HEM	C3C-C2C	-2.55	1.39	1.43
10	N	1304	TDS	CAR-CAQ	-2.54	1.39	1.52
9	A	301	HEM	CHC-C1C	2.53	1.40	1.36
9	N	301	HEM	FE-NC	2.53	2.07	1.97
11	Q	1305	PL9	C7-C3	2.53	1.53	1.51
13	B	201	CLA	C1B-C2B	2.52	1.43	1.40
9	A	302	HEM	C3D-C2D	-2.51	1.39	1.43
10	A	304	TDS	CAR-CAQ	-2.51	1.39	1.52
9	A	302	HEM	FE-NC	2.51	2.07	1.97
9	N	303	HEM	CHC-C1C	2.50	1.40	1.36
13	O	1201	CLA	MG-ND	-2.49	1.99	2.05
9	A	301	HEM	C1A-C2A	2.49	1.47	1.43
12	Q	1307	OPC	CBQ-CBR	-2.49	1.39	1.50
9	A	303	HEM	CMD-C2D	2.47	1.55	1.47
12	N	1306	OPC	CBT-CBS	-2.47	1.39	1.50
12	Q	1307	OPC	CBT-CBS	-2.47	1.39	1.50
12	N	1306	OPC	CBQ-CBR	-2.46	1.39	1.50
12	B	307	OPC	CBT-CBS	-2.46	1.39	1.50
9	N	303	HEM	CHD-C4C	2.46	1.40	1.36
9	P	301	HEM	C3B-CAB	2.46	1.48	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
13	B	201	CLA	C2A-C1A	-2.45	1.47	1.52
12	D	306	OPC	CBT-CBS	-2.45	1.39	1.50
12	B	307	OPC	CBQ-CBR	-2.44	1.39	1.50
9	A	301	HEM	FE-NC	2.44	2.07	1.97
12	D	306	OPC	CBQ-CBR	-2.43	1.39	1.50
13	O	1201	CLA	C1B-NB	2.43	1.37	1.34
9	N	303	HEM	CAA-C2A	2.42	1.56	1.52
9	N	302	HEM	C3B-CAB	2.39	1.47	1.40
9	N	301	HEM	CHA-C4D	2.39	1.39	1.35
9	N	302	HEM	CMA-C3A	2.33	1.56	1.51
13	B	201	CLA	OBD-CAD	2.28	1.25	1.22
9	C	301	HEM	CHC-C1C	2.28	1.40	1.36
9	N	303	HEM	FE-NA	2.25	2.02	1.92
9	A	302	HEM	FE-NA	2.25	2.02	1.92
9	A	303	HEM	CMB-C2B	2.24	1.54	1.47
9	A	302	HEM	CMA-C3A	2.18	1.56	1.51
15	E	101	BCR	C8-C9	2.17	1.50	1.45
9	P	301	HEM	C3D-C2D	-2.15	1.40	1.43
9	N	303	HEM	C4A-NA	2.14	1.40	1.36
9	N	303	HEM	CMB-C2B	2.13	1.54	1.47
9	N	302	HEM	CAA-C2A	-2.12	1.48	1.52
9	A	303	HEM	CHC-C1C	2.11	1.40	1.36
9	N	301	HEM	CMA-C3A	2.11	1.56	1.51
9	P	301	HEM	CBA-CGA	2.11	1.55	1.50
9	C	301	HEM	C1A-C2A	2.09	1.47	1.43
9	A	301	HEM	CMA-C3A	2.08	1.56	1.51
13	B	201	CLA	CBD-CHA	2.07	1.60	1.53
13	B	201	CLA	CBA-CGA	2.06	1.56	1.50
15	R	1101	BCR	C14-C13	2.05	1.38	1.35
13	B	201	CLA	MG-ND	-2.05	2.00	2.05
9	C	301	HEM	C3D-C2D	-2.02	1.40	1.43
9	P	301	HEM	CHB-C1B	2.01	1.38	1.35
13	O	1201	CLA	C2A-C1A	-2.01	1.48	1.52
9	C	301	HEM	C4D-ND	-2.00	1.35	1.39
13	O	1201	CLA	CAC-C3C	2.00	1.56	1.51
13	B	201	CLA	C3B-C2B	-2.00	1.37	1.41

All (262) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	A	304	TDS	CAP-OAO-CAN	-21.65	115.01	122.07
10	N	1304	TDS	CAP-OAO-CAN	-18.60	116.01	122.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	N	302	HEM	CBA-CAA-C2A	-15.21	85.90	112.69
9	A	302	HEM	CBA-CAA-C2A	-14.90	86.43	112.69
10	N	1304	TDS	OAO-CAP-CAQ	11.80	117.17	110.58
10	A	304	TDS	OAO-CAP-CAQ	10.23	116.29	110.58
13	O	1201	CLA	CAA-C2A-C1A	8.09	131.82	111.62
11	A	305	PL9	C7-C8-C9	-8.01	113.23	126.76
11	Q	1305	PL9	C7-C8-C9	-7.94	113.34	126.76
13	B	201	CLA	CAA-C2A-C1A	7.80	131.10	111.62
10	A	304	TDS	CAL-CAM-CAN	-7.56	112.94	120.27
11	Q	1305	PL9	C17-C18-C19	-7.52	111.57	127.80
10	N	1304	TDS	CAL-CAM-CAN	-7.42	113.07	120.27
13	O	1201	CLA	C4D-C3D-CAD	7.36	117.10	108.05
11	A	305	PL9	C22-C23-C24	-7.28	112.11	127.80
11	Q	1305	PL9	C12-C13-C14	-7.21	112.24	127.80
11	Q	1305	PL9	C37-C38-C39	-7.04	112.62	127.80
11	A	305	PL9	C17-C18-C19	-7.00	112.69	127.80
15	R	1101	BCR	C38-C26-C25	6.94	132.37	124.51
11	A	305	PL9	C37-C38-C39	-6.93	112.84	127.80
13	B	201	CLA	CMA-C3A-C4A	6.90	131.78	111.76
9	A	303	HEM	C2D-C1D-ND	-6.83	104.86	112.93
11	A	305	PL9	C12-C13-C14	-6.79	113.15	127.80
13	B	201	CLA	C4D-C3D-CAD	6.76	116.36	108.05
11	Q	1305	PL9	C22-C23-C24	-6.72	113.31	127.80
9	N	303	HEM	C3B-C4B-NB	-6.62	109.27	114.00
15	E	101	BCR	C38-C26-C25	6.56	131.94	124.51
13	O	1201	CLA	CMA-C3A-C4A	6.45	130.49	111.76
9	A	303	HEM	C3B-C4B-NB	-6.44	109.39	114.00
9	A	302	HEM	CAA-CBA-CGA	6.16	133.27	113.47
11	Q	1305	PL9	C25-C24-C26	5.77	124.16	115.39
9	N	302	HEM	CAA-CBA-CGA	5.75	131.95	113.47
9	A	301	HEM	C3B-C4B-NB	-5.68	109.93	114.00
10	N	1304	TDS	CAE-CAF-CAN	5.62	122.41	117.08
9	N	303	HEM	C2D-C1D-ND	-5.59	106.33	112.93
15	R	1101	BCR	C24-C23-C22	5.57	134.54	126.22
9	N	301	HEM	C3B-C4B-NB	-5.49	110.07	114.00
10	A	304	TDS	CAJ-OAK-CAL	-5.35	109.67	117.59
9	A	302	HEM	C3B-C4B-NB	-5.21	110.27	114.00
10	N	1304	TDS	CAJ-OAK-CAL	-5.19	109.90	117.59
15	E	101	BCR	C24-C23-C22	5.19	133.98	126.22
10	A	304	TDS	CAE-CAF-CAN	5.12	121.94	117.08
13	B	201	CLA	C3D-CAD-CBD	-5.12	100.36	107.60
15	E	101	BCR	C20-C21-C22	4.93	134.39	127.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	C	301	HEM	C3B-C4B-NB	-4.91	110.49	114.00
10	A	304	TDS	CAA-OAB-CAE	-4.89	111.22	117.78
15	E	101	BCR	C33-C5-C6	4.86	130.01	124.51
9	N	302	HEM	C3B-C4B-NB	-4.85	110.53	114.00
13	B	201	CLA	CBC-CAC-C3C	4.84	127.01	112.38
15	R	1101	BCR	C33-C5-C6	4.83	129.98	124.51
13	O	1201	CLA	C3D-CAD-CBD	-4.80	100.81	107.60
15	R	1101	BCR	C20-C21-C22	4.79	134.20	127.29
12	N	1306	OPC	OAN-CAO-CAP	4.75	125.24	110.51
9	P	301	HEM	C3B-C4B-NB	-4.74	110.61	114.00
11	A	305	PL9	C25-C24-C26	4.62	122.42	115.39
13	O	1201	CLA	C3A-C2A-C1A	4.62	107.61	101.08
13	O	1201	CLA	CBD-CHA-C1A	4.38	134.50	128.77
13	O	1201	CLA	C2A-C1A-NA	-4.38	106.40	111.24
9	A	303	HEM	CAD-C3D-C4D	4.36	132.37	124.53
10	N	1304	TDS	CAA-OAB-CAE	-4.36	111.93	117.78
13	B	201	CLA	C1D-C2D-C3D	-4.31	103.26	106.78
13	O	1201	CLA	C1D-C2D-C3D	-4.25	103.31	106.78
13	O	1201	CLA	CBC-CAC-C3C	4.23	125.17	112.38
12	Q	1307	OPC	OAN-CAO-CAP	4.21	123.56	110.51
13	B	201	CLA	C3A-C2A-C1A	4.21	107.03	101.08
9	A	303	HEM	CBA-CAA-C2A	-4.17	105.35	112.69
13	B	201	CLA	C2A-C3A-C4A	4.13	107.75	101.40
15	R	1101	BCR	C38-C26-C27	-4.12	105.77	113.34
15	E	101	BCR	C7-C8-C9	4.05	132.28	126.22
12	D	306	OPC	OAN-CAO-CAP	4.02	122.96	110.51
13	O	1201	CLA	C2A-C3A-C4A	3.96	107.48	101.40
15	E	101	BCR	C38-C26-C27	-3.95	106.08	113.34
15	R	1101	BCR	C23-C22-C21	-3.93	112.93	118.97
10	N	1304	TDS	OAB-CAE-CAD	-3.84	117.78	123.59
11	Q	1305	PL9	C15-C14-C16	3.81	121.18	115.39
10	A	304	TDS	CAQ-CAP-CAH	3.80	125.01	120.42
15	R	1101	BCR	C33-C5-C4	-3.80	106.37	113.34
13	O	1201	CLA	C2A-C1A-CHA	3.80	130.41	123.83
15	R	1101	BCR	C7-C8-C9	3.79	131.89	126.22
15	E	101	BCR	C33-C5-C4	-3.76	106.43	113.34
15	R	1101	BCR	C30-C25-C26	-3.75	117.18	122.60
11	Q	1305	PL9	C26-C24-C23	-3.74	113.89	121.08
15	E	101	BCR	C30-C25-C26	-3.73	117.20	122.60
15	E	101	BCR	C23-C22-C21	-3.73	113.24	118.97
12	B	307	OPC	OAN-CAO-CAP	3.71	122.02	110.51
13	B	201	CLA	OBD-CAD-C3D	3.71	134.81	127.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	N	1306	OPC	OBJ-CBK-CBL	3.69	123.55	111.94
11	Q	1305	PL9	C30-C29-C31	3.62	120.90	115.39
9	N	301	HEM	CBA-CAA-C2A	-3.60	106.35	112.69
13	O	1201	CLA	C4A-NA-C1A	3.58	111.46	106.52
15	R	1101	BCR	C29-C30-C25	3.56	116.35	110.44
13	B	201	CLA	C1-C2-C3	3.53	132.46	126.19
15	E	101	BCR	C29-C30-C25	3.51	116.28	110.44
9	N	303	HEM	CBD-CAD-C3D	-3.48	106.78	114.37
13	B	201	CLA	C3A-C4A-NA	-3.38	106.83	110.95
11	A	305	PL9	C30-C29-C31	3.36	120.50	115.39
9	N	301	HEM	CBD-CAD-C3D	-3.36	107.05	114.37
13	O	1201	CLA	C1-C2-C3	3.35	132.14	126.19
10	N	1304	TDS	CAQ-CAP-CAH	3.34	124.45	120.42
13	B	201	CLA	C2A-C1A-NA	-3.29	107.60	111.24
13	O	1201	CLA	OBD-CAD-C3D	3.29	134.03	127.91
13	O	1201	CLA	C3A-C4A-NA	-3.21	107.03	110.95
15	E	101	BCR	C1-C6-C5	-3.21	117.96	122.60
9	A	303	HEM	CAD-CBD-CGD	3.20	123.46	113.48
13	B	201	CLA	CMD-C2D-C3D	3.18	129.98	124.97
12	D	306	OPC	OBJ-CBK-CBL	3.18	121.94	111.94
9	N	301	HEM	C2D-C1D-ND	-3.17	109.19	112.93
9	A	302	HEM	C2D-C1D-ND	-3.16	109.19	112.93
15	E	101	BCR	C2-C1-C6	3.16	115.70	110.44
9	P	301	HEM	CAA-C2A-C3A	-3.14	120.03	129.00
15	R	1101	BCR	C1-C6-C5	-3.14	118.05	122.60
10	A	304	TDS	CAG-CAF-CAE	-3.13	120.42	125.23
13	B	201	CLA	CMB-C2B-C1B	-3.11	123.84	128.62
15	R	1101	BCR	C2-C1-C6	3.10	115.59	110.44
11	Q	1305	PL9	C35-C34-C36	3.09	120.09	115.39
15	R	1101	BCR	C11-C10-C9	3.08	131.73	127.29
13	B	201	CLA	C4A-NA-C1A	3.06	110.73	106.52
11	A	305	PL9	C26-C24-C23	-3.05	115.21	121.08
9	C	301	HEM	CAA-C2A-C3A	-3.04	120.31	129.00
9	N	302	HEM	CAA-C2A-C3A	-3.00	120.43	129.00
10	N	1304	TDS	CAG-CAF-CAE	-3.00	120.61	125.23
10	A	304	TDS	OAB-CAE-CAD	-3.00	119.06	123.59
15	E	101	BCR	C23-C24-C25	3.00	136.18	127.32
13	B	201	CLA	C2A-C1A-CHA	2.99	129.00	123.83
9	A	301	HEM	C2D-C1D-ND	-2.97	109.42	112.93
13	O	1201	CLA	C4D-CHA-CBD	-2.96	102.39	109.37
13	B	201	CLA	CBD-CHA-C1A	2.96	132.64	128.77
13	O	1201	CLA	CMB-C2B-C1B	-2.96	124.07	128.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	N	302	HEM	C2D-C1D-ND	-2.90	109.50	112.93
15	R	1101	BCR	C23-C24-C25	2.90	135.88	127.32
9	N	301	HEM	CAA-C2A-C3A	-2.89	120.74	129.00
13	O	1201	CLA	CAA-C2A-C3A	2.89	119.88	113.04
9	A	303	HEM	CAA-CBA-CGA	2.89	122.75	113.47
11	A	305	PL9	C10-C9-C8	-2.87	117.84	123.52
12	B	307	OPC	CBQ-CBR-CBS	-2.85	109.28	125.43
9	A	301	HEM	CAA-C2A-C3A	-2.83	120.92	129.00
9	A	301	HEM	CHA-C4D-ND	2.81	128.17	124.31
9	N	301	HEM	CHA-C4D-ND	2.79	128.15	124.31
9	N	303	HEM	C2A-C1A-NA	2.78	113.60	109.73
9	P	301	HEM	C2D-C1D-ND	-2.78	109.65	112.93
15	E	101	BCR	C1-C6-C7	2.75	123.32	115.69
9	N	303	HEM	C4A-NA-C1A	-2.73	103.17	106.76
15	R	1101	BCR	C1-C6-C7	2.73	123.24	115.69
11	A	305	PL9	C35-C34-C36	2.72	119.53	115.39
9	A	303	HEM	C1A-CHA-C4D	-2.72	123.90	127.47
15	E	101	BCR	C15-C14-C13	2.71	131.20	127.29
15	E	101	BCR	C11-C10-C9	2.69	131.17	127.29
9	N	302	HEM	C1A-C2A-C3A	2.69	109.70	106.92
11	Q	1305	PL9	C42-C43-C44	2.68	133.57	127.80
13	B	201	CLA	CAA-C2A-C3A	2.66	119.33	113.04
11	Q	1305	PL9	C10-C9-C11	2.66	119.43	115.39
9	A	303	HEM	C2A-C1A-NA	2.65	113.41	109.73
9	A	303	HEM	C4A-NA-C1A	-2.64	103.28	106.76
9	N	302	HEM	C4A-NA-C1A	2.64	110.24	106.76
13	O	1201	CLA	CMD-C2D-C3D	2.64	129.12	124.97
12	N	1306	OPC	OAN-CAO-OAD	-2.63	117.06	122.95
12	D	306	OPC	CAX-CAW-CAV	-2.63	110.52	125.43
11	Q	1305	PL9	C20-C19-C21	2.62	119.37	115.39
15	E	101	BCR	C35-C13-C12	2.62	122.32	118.09
9	N	301	HEM	C4C-NC-C1C	2.62	108.26	105.53
11	A	305	PL9	C20-C19-C21	2.61	119.36	115.39
13	B	201	CLA	O2A-CGA-CBA	2.61	120.15	111.94
12	B	307	OPC	CAX-CAW-CAV	-2.61	110.63	125.43
12	Q	1307	OPC	OAN-CAO-OAD	-2.60	117.13	122.95
11	A	305	PL9	C45-C44-C43	-2.60	118.38	123.52
11	A	305	PL9	C42-C43-C44	2.59	133.39	127.80
15	R	1101	BCR	C15-C14-C13	2.58	131.01	127.29
12	B	307	OPC	CBM-CBL-CBK	2.58	123.62	113.51
13	B	201	CLA	O2D-CGD-CBD	2.57	116.57	111.33
12	Q	1307	OPC	OBJ-CBK-CBL	2.57	120.02	111.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	R	1101	BCR	C35-C13-C12	2.55	122.21	118.09
15	E	101	BCR	C37-C22-C23	2.54	122.20	118.09
15	R	1101	BCR	C37-C22-C23	2.54	122.19	118.09
9	C	301	HEM	C2D-C1D-ND	-2.53	109.95	112.93
9	A	302	HEM	C1A-C2A-C3A	2.52	109.53	106.92
12	N	1306	OPC	CBQ-CBR-CBS	-2.52	111.14	125.43
11	Q	1305	PL9	C40-C39-C41	2.52	119.22	115.39
11	A	305	PL9	C40-C39-C41	2.52	119.22	115.39
13	O	1201	CLA	O2A-CGA-CBA	2.51	119.85	111.94
15	E	101	BCR	C12-C13-C14	-2.51	115.12	118.97
13	O	1201	CLA	C11-C12-C13	2.50	122.34	115.14
9	A	301	HEM	C4C-NC-C1C	2.50	108.13	105.53
15	R	1101	BCR	C12-C13-C14	-2.49	115.14	118.97
9	A	302	HEM	CAA-C2A-C3A	-2.48	121.92	129.00
11	Q	1305	PL9	C10-C9-C8	-2.48	118.61	123.52
9	A	302	HEM	CMA-C3A-C4A	2.48	132.43	128.62
9	A	301	HEM	CBD-CAD-C3D	-2.48	108.97	114.37
15	E	101	BCR	C32-C1-C6	2.47	114.42	110.33
9	N	302	HEM	C2A-C1A-NA	-2.45	106.33	109.73
12	Q	1307	OPC	CBQ-CBR-CBS	-2.45	111.55	125.43
15	E	101	BCR	C34-C9-C10	-2.45	119.44	122.92
9	P	301	HEM	CHA-C4D-ND	2.44	127.66	124.31
10	N	1304	TDS	OAB-CAE-CAF	2.43	120.63	116.59
9	C	301	HEM	C1A-CHA-C4D	-2.42	124.29	127.47
12	N	1306	OPC	CAX-CAW-CAV	-2.42	111.74	125.43
9	N	302	HEM	C3A-C4A-NA	-2.42	107.59	109.41
15	R	1101	BCR	C8-C7-C6	2.40	134.41	127.32
11	Q	1305	PL9	C40-C39-C38	-2.38	118.81	123.52
11	A	305	PL9	C10-C9-C11	2.37	119.00	115.39
12	Q	1307	OPC	CBI-OBJ-CBK	-2.37	110.64	116.04
13	B	201	CLA	C4B-C3B-CAB	2.37	131.97	127.18
9	A	302	HEM	C4C-NC-C1C	2.36	107.99	105.53
11	Q	1305	PL9	C30-C29-C28	-2.36	118.85	123.52
11	Q	1305	PL9	C45-C44-C43	-2.35	118.87	123.52
11	Q	1305	PL9	C35-C34-C33	-2.34	118.89	123.52
9	P	301	HEM	CAA-C2A-C1A	2.33	132.53	125.50
9	A	303	HEM	CHD-C4C-NC	-2.33	122.71	124.73
15	E	101	BCR	C8-C7-C6	2.32	134.18	127.32
12	D	306	OPC	OAN-CAO-OAD	-2.32	117.75	122.95
11	A	305	PL9	C40-C39-C38	-2.32	118.92	123.52
10	N	1304	TDS	OAO-CAN-CAM	2.31	118.88	115.90
9	N	303	HEM	CMA-C3A-C4A	-2.31	125.06	128.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	P	301	HEM	C1A-CHA-C4D	-2.31	124.44	127.47
9	C	301	HEM	CHA-C4D-ND	2.30	127.47	124.31
15	R	1101	BCR	C32-C1-C6	2.29	114.12	110.33
13	O	1201	CLA	CED-O2D-CGD	2.29	121.46	116.02
12	N	1306	OPC	CBO-CBP-CBQ	2.28	122.46	113.78
9	N	303	HEM	C4D-ND-C1D	2.28	107.50	105.16
9	A	303	HEM	C4D-ND-C1D	2.24	107.45	105.16
9	A	302	HEM	C4A-NA-C1A	2.23	109.70	106.76
9	N	303	HEM	CAD-CBD-CGD	2.23	120.43	113.48
9	A	301	HEM	CAA-C2A-C1A	2.23	132.22	125.50
13	B	201	CLA	C4D-CHA-CBD	-2.21	104.16	109.37
9	C	301	HEM	C3A-C4A-NA	-2.21	107.75	109.41
9	N	301	HEM	CAA-C2A-C1A	2.20	132.15	125.50
9	A	303	HEM	CHD-C1D-ND	2.19	126.41	124.58
11	A	305	PL9	C53-C6-C1	2.19	119.58	114.80
11	A	305	PL9	C15-C14-C13	-2.19	119.19	123.52
11	Q	1305	PL9	C20-C19-C18	-2.19	119.19	123.52
9	C	301	HEM	CAA-C2A-C1A	2.17	132.05	125.50
9	C	301	HEM	C4A-NA-C1A	2.16	109.61	106.76
15	E	101	BCR	C34-C9-C8	2.16	121.58	118.09
11	A	305	PL9	C7-C3-C4	-2.16	114.96	117.10
11	A	305	PL9	C45-C44-C46	2.15	118.66	115.39
11	Q	1305	PL9	C45-C44-C46	2.15	118.66	115.39
10	A	304	TDS	OBD-CAM-CAN	2.14	123.67	119.81
9	A	301	HEM	CMA-C3A-C2A	-2.14	120.91	124.94
11	Q	1305	PL9	C53-C6-C1	2.13	119.46	114.80
10	A	304	TDS	CAH-CAG-CAF	-2.12	118.34	121.65
12	N	1306	OPC	OBJ-CBK-OCC	-2.12	117.65	123.43
13	B	201	CLA	C4B-C3B-C2B	-2.12	104.41	106.97
12	N	1306	OPC	CBM-CBL-CBK	2.11	121.80	113.51
9	P	301	HEM	CMA-C3A-C2A	-2.11	120.97	124.94
9	N	302	HEM	C4C-NC-C1C	2.10	107.72	105.53
11	A	305	PL9	C35-C34-C33	-2.10	119.37	123.52
10	N	1304	TDS	OAO-CAP-CAH	-2.09	118.31	120.92
10	N	1304	TDS	OBD-CAM-CAN	2.09	123.57	119.81
12	B	307	OPC	CBO-CBP-CBQ	2.09	121.71	113.78
9	A	302	HEM	C3A-C4A-NA	-2.08	107.84	109.41
9	N	302	HEM	CMA-C3A-C4A	2.08	131.82	128.62
9	A	303	HEM	CAA-C2A-C3A	-2.07	123.09	129.00
12	B	307	OPC	OAN-CAO-OAD	-2.06	118.34	122.95
15	R	1101	BCR	C34-C9-C8	2.06	121.42	118.09
9	N	301	HEM	CAA-CBA-CGA	2.05	120.07	113.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	301	HEM	CMA-C3A-C4A	2.05	131.78	128.62
9	A	302	HEM	O2A-CGA-CBA	2.04	121.43	114.22
10	A	304	TDS	OAO-CAN-CAM	2.03	118.52	115.90
9	N	302	HEM	C1A-CHA-C4D	-2.02	124.81	127.47
9	A	303	HEM	C3A-C4A-NA	2.01	110.93	109.41
9	A	301	HEM	C1B-NB-C4B	2.01	107.22	105.16
9	C	301	HEM	CAD-C3D-C4D	2.01	128.13	124.53
9	N	302	HEM	O1A-CGA-CBA	-2.00	116.13	123.03

All (5) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
13	B	201	CLA	C8
13	B	201	CLA	C3A
13	O	1201	CLA	C8
13	O	1201	CLA	C2A
13	O	1201	CLA	C3A

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
12	Q	1307	OPC	CAM-OAN-CAO-CAP
12	Q	1307	OPC	CAM-OAN-CAO-OAD
12	B	307	OPC	CAM-OAN-CAO-OAD
12	B	307	OPC	CAM-OAN-CAO-CAP
12	D	306	OPC	CAM-OAN-CAO-CAP
12	N	1306	OPC	CAM-OAN-CAO-CAP
12	D	306	OPC	CAM-OAN-CAO-OAD
12	N	1306	OPC	CAM-OAN-CAO-OAD
11	A	305	PL9	C44-C43-C42-C41
11	Q	1305	PL9	C44-C43-C42-C41
12	D	306	OPC	CAO-OAN-CAM-CAL
12	N	1306	OPC	CAO-OAN-CAM-CAL

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	202/215 (93%)	-0.13	0	100	100	22, 55, 101, 174	0
1	N	202/215 (93%)	-0.14	0	100	100	19, 53, 86, 121	0
2	B	138/160 (86%)	-0.16	2 (1%)	72	18	24, 65, 122, 194	0
2	O	138/160 (86%)	-0.11	0	100	100	28, 62, 139, 184	0
3	C	286/289 (98%)	-0.08	2 (0%)	84	28	5, 82, 153, 200	1 (0%)
3	P	286/289 (98%)	-0.02	2 (0%)	84	28	14, 85, 154, 200	1 (0%)
4	D	168/179 (93%)	-0.04	2 (1%)	75	20	36, 117, 179, 200	0
4	Q	168/179 (93%)	-0.09	0	100	100	33, 109, 165, 192	0
5	E	32/32 (100%)	0.10	0	100	100	30, 76, 149, 177	0
5	R	32/32 (100%)	0.13	0	100	100	37, 81, 149, 171	0
6	F	33/35 (94%)	-0.08	0	100	100	39, 63, 140, 177	0
6	S	35/35 (100%)	-0.13	0	100	100	41, 78, 152, 167	0
7	G	23/37 (62%)	0.01	0	100	100	33, 71, 128, 171	0
7	T	27/37 (72%)	-0.07	0	100	100	37, 65, 100, 152	0
8	H	27/29 (93%)	-0.11	0	100	100	33, 71, 130, 165	0
8	U	27/29 (93%)	-0.08	0	100	100	46, 81, 152, 157	0
All	All	1824/1952 (93%)	-0.08	8 (0%)	90	41	5, 74, 155, 200	2 (0%)

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	117	GLY	3.2
2	B	93	ASN	2.5
4	D	16	ARG	2.4
3	P	101	VAL	2.3
3	P	71	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
3	C	73	GLY	2.1
4	D	91	ILE	2.1
2	B	95	LEU	2.1

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

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

## 6.3 Carbohydrates ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
10	TDS	N	1304	30/30	0.45	3.75	80,80,80,80	0
11	PL9	A	305	55/55	0.41	3.47	98,98,98,98	0
15	BCR	R	1101	40/40	0.56	3.46	91,91,91,91	0
15	BCR	E	101	40/40	0.47	3.16	77,77,77,77	0
9	HEM	A	301	43/43	0.28	2.33	45,45,45,45	0
13	CLA	B	201	65/65	0.28	2.19	45,56,56,56	0
9	HEM	N	302	43/43	0.28	1.91	48,56,56,56	0
12	OPC	N	1306	54/55	0.32	1.87	87,87,87,87	0
9	HEM	N	301	43/43	0.28	1.83	41,50,50,50	0
9	HEM	A	302	43/43	0.29	1.67	62,62,62,62	0
11	PL9	Q	1305	55/55	0.29	1.31	85,85,85,85	0
10	TDS	A	304	30/30	0.34	1.26	52,52,52,52	0
13	CLA	O	1201	65/65	0.27	1.05	22,72,72,72	0
12	OPC	D	306	54/55	0.29	1.04	86,86,86,86	0
9	HEM	A	303	43/43	0.25	0.40	55,81,81,81	0
9	HEM	C	301	43/43	0.26	0.11	62,83,83,83	0
12	OPC	B	307	54/55	0.24	0.10	84,84,84,84	0
9	HEM	N	303	43/43	0.22	-0.19	61,61,61,64	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
9	HEM	P	301	43/43	0.24	-0.26	47,56,56,56	0
12	OPC	Q	1307	54/55	0.20	-0.69	72,72,72,72	0
14	FES	Q	1200	4/4	0.15	-0.92	94,94,98,98	0
14	FES	D	200	4/4	0.10	-1.56	144,144,170,170	0

## 6.5 Other polymers

There are no such residues in this entry.