



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 09:16 AM GMT

PDB ID : 1A3X
Title : PYRUVATE KINASE FROM SACCHAROMYCES CEREVISIAE COM-
PLEXED WITH PG, MN²⁺ AND K⁺
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Deposited on : 1998-01-26
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
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

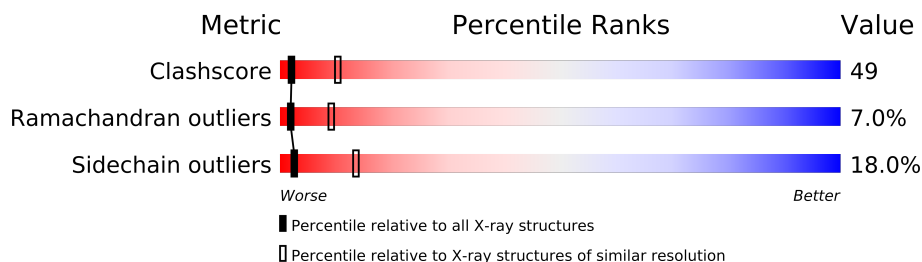
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
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))
Clashscore	79885	1594 (3.00-3.00)
Ramachandran outliers	78287	1537 (3.00-3.00)
Sidechain outliers	78261	1540 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	500	
1	B	500	

2 Entry composition i

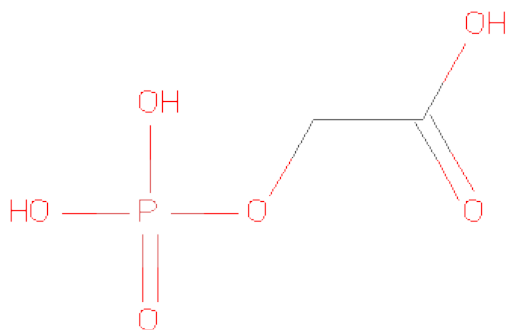
There are 4 unique types of molecules in this entry. The entry contains 7472 atoms, of which 0 are hydrogen and 0 are deuterium.

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 PYRUVATE KINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	487	Total	C	N	O	S	0	0	0
			3725	2347	642	718	18			
1	B	487	Total	C	N	O	S	0	0	0
			3725	2347	642	718	18			

- Molecule 2 is SUGAR (2-PHOSPHOGLYCOLIC ACID) (three-letter code: PGA) (formula: $C_2H_5O_6P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			9	2	6	1		
2	B	1	Total	C	O	P	0	0
			9	2	6	1		

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Mn 1	0	0
3	A	1	Total 1	Mn 1	0	0

- Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	K 1	0	0
4	A	1	Total 1	K 1	0	0

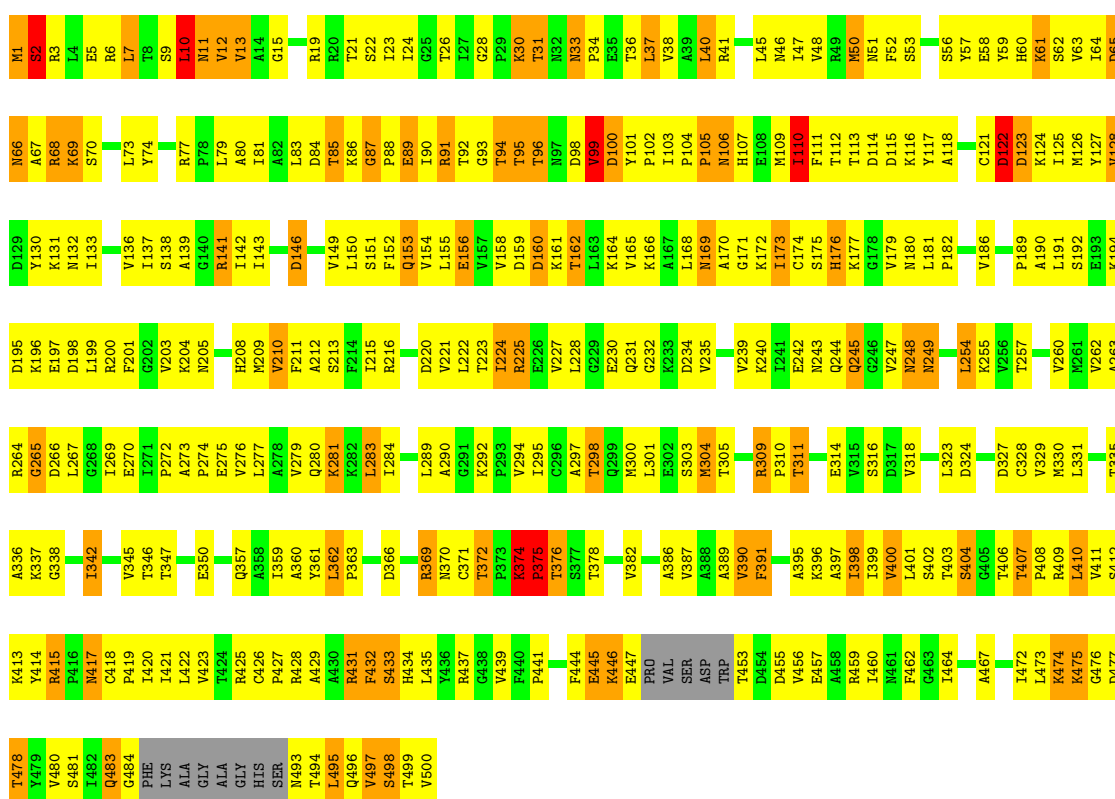
3 Residue-property plots

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.

Note EDS was not executed.

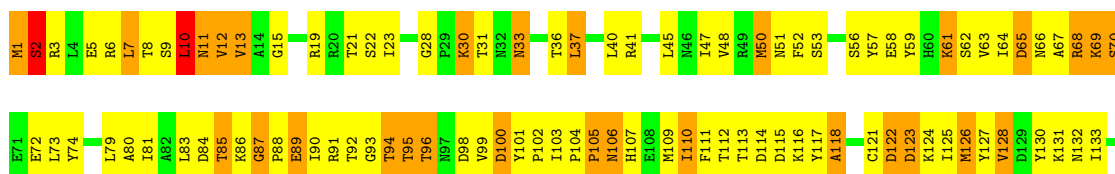
• Molecule 1: PYRUVATE KINASE

Chain A:



• Molecule 1: PYRUVATE KINASE

Chain B:



PHE	L422	M348	L273	E197	V136
LYS	V423	A349	P272	D198	I137
ALA	T425	E350	P273	L199	S138
GLY	R425	Q357	P274	R200	A139
ALA	C426	A358	E275	F201	G140
GLY	P427	I359			R141
HIS	R428	A360	V279	H208	I142
SER	A429	Y361	Q280	M209	I143
N493	A430	K362	K281	V210	Y144
T494	R431	L362	L282	F211	V145
L495	F432	P363	L283	A212	D146
Q496	S433	D366	L284	S213	D147
H434	H434		L289	F214	G148
V497	L435	R369		I215	V149
S498	T436	N370	K292	R216	L150
V500	R437	T371	P293	D220	S151
	G438	C372	V294	D221	F152
	V439	T372	P294	V221	I153
	F440	P373	L295	L222	V154
	P441	K374	C296	T223	L155
		P375	A297	I224	E156
	F444	T376	T298	E226	V157
E445	E445	S377	Q299	R225	V158
K446	K446	T378	N300	V227	D159
E447	E447	L301	L301	L228	D160
PRO	PRO	V382	E302		K161
VAL	VAL		S303	Q231	T162
SER	SER	A386	K304	G232	L163
ASP	ASP	V387	T305		K164
TRP	TRP	A388		V235	V165
T453	T453	A389	P308		V166
D454	D454	V390	R309	V239	A167
V456	V456	F391	P310	K240	L168
E457	E457		T311	I241	N169
A458	A458	A395	E314	E242	A170
R459	R459	K396	V315	N243	G171
I460	I460	I398	S316	Q244	K172
R461	R461	I399	D317	Q245	I173
F462	F462	V400	V318	G246	C174
G463	G463	L401		V247	S175
I464	I464	S402	L323	N248	H176
E465	E465	T403	D324	N249	K177
K466	K466	S404			G178
K467	K467		D327	L254	V179
F468	F468	T407	C328	K255	N180
E469	E469	P408	V329	V256	L181
		R409	L330	D258	P182
		L410	L331	G259	G183
L473	L473	V411		V260	T184
K474	K474	S412	A336	D261	D185
K475	K475	K413	K337	V262	V186
G476	G476	Y414	G338	A263	D187
D477	D477	R415		R264	L188
T478	T478	P416	I342	G265	P189
V479	V479	N417	N343	D266	A190
V480	V480	C418	K344	L267	L191
S481	S481	P419	V345	G268	
T482	T482	I420	T346	I269	K194
Q483	Q483	I421	T347	E270	D195
G484	G484				K196

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	108.30Å 106.40Å 105.50Å 90.00° 110.80° 90.00°	Depositor
Resolution (Å)	100.00 – 3.00	Depositor
% Data completeness (in resolution range)	85.0 (100.00-3.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.227 , 0.341	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7472	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: K, MN, PGA

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	0.59	2/3781 (0.1%)	0.88	4/5124 (0.1%)
1	B	0.41	0/3781	1.03	7/5124 (0.1%)
All	All	0.51	2/7562 (0.0%)	0.96	11/10248 (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	A	0	2
1	B	0	2
All	All	0	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	374	LYS	C-N	-26.09	0.84	1.34
1	A	375	PRO	C-N	5.31	1.46	1.34

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	375	PRO	O-C-N	-31.91	71.64	122.70
1	B	375	PRO	CA-C-N	-31.09	48.81	117.20
1	A	375	PRO	O-C-N	-25.38	82.10	122.70
1	A	375	PRO	CA-C-N	-24.09	64.20	117.20
1	B	374	LYS	O-C-N	-21.78	79.71	121.10
1	B	374	LYS	C-N-CD	-21.33	73.68	120.60
1	A	374	LYS	C-N-CD	-20.82	74.79	120.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	374	LYS	CA-C-N	14.60	157.98	117.10
1	B	375	PRO	C-N-CA	-7.56	102.79	121.70
1	A	374	LYS	C-N-CA	-7.31	91.29	122.00
1	B	374	LYS	C-N-CA	-5.03	100.89	122.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	374	LYS	Peptide
1	A	375	PRO	Mainchain
1	B	374	LYS	Mainchain,Peptide

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	3725	0	3804	369	1
1	B	3725	0	3804	375	1
2	A	9	0	2	5	0
2	B	9	0	2	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
All	All	7472	0	7612	744	1

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 49.

All (744) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:374:LYS:HG2	1:A:375:PRO:HD3	1.25	1.11
1:B:186:VAL:HG23	1:B:216:ARG:HE	1.19	1.06
1:A:272:PRO:HB2	1:A:275:GLU:HG3	1.28	1.06

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:272:PRO:HB2	1:B:275:GLU:HG3	1.37	1.05
1:A:390:VAL:HB	1:A:395:ALA:HB3	1.37	1.02
1:A:378:THR:HG22	1:A:493:ASN:HD21	1.18	1.01
1:A:263:ALA:HB1	2:A:1005:PGA:O2	1.58	1.00
1:A:374:LYS:HG2	1:A:375:PRO:CD	1.89	1.00
1:B:378:THR:HG22	1:B:493:ASN:HD21	1.21	0.98
1:B:375:PRO:HB2	1:B:376:THR:OG1	1.61	0.98
1:B:390:VAL:HB	1:B:395:ALA:HB3	1.45	0.96
1:A:263:ALA:CB	2:A:1005:PGA:O2	2.14	0.96
1:A:220:ASP:O	1:A:224:ILE:HG22	1.63	0.95
1:B:374:LYS:H	1:B:375:PRO:HB3	1.31	0.95
1:B:374:LYS:N	1:B:375:PRO:HB3	1.82	0.94
1:B:220:ASP:O	1:B:224:ILE:HG22	1.69	0.93
1:B:342:ILE:HD13	1:B:342:ILE:H	1.33	0.91
1:A:374:LYS:CG	1:A:375:PRO:HD3	2.03	0.89
1:B:375:PRO:CB	1:B:376:THR:OG1	2.19	0.89
1:A:342:ILE:HD13	1:A:342:ILE:H	1.36	0.89
1:A:369:ARG:HE	1:A:370:ASN:HD21	1.22	0.87
1:B:310:PRO:HG3	1:B:347:THR:HG21	1.58	0.86
1:B:395:ALA:HA	1:B:478:THR:HG23	1.57	0.84
1:A:264:ARG:HH21	1:A:280:GLN:HE22	1.24	0.84
1:A:224:ILE:O	1:A:228:LEU:HD23	1.78	0.84
1:A:265:GLY:N	2:A:1005:PGA:O1	2.12	0.83
1:A:378:THR:HG22	1:A:493:ASN:ND2	1.93	0.83
1:A:395:ALA:HA	1:A:478:THR:HG23	1.61	0.83
1:B:224:ILE:O	1:B:228:LEU:HD23	1.79	0.83
1:A:310:PRO:HG3	1:A:347:THR:HG21	1.61	0.83
1:B:378:THR:HG22	1:B:493:ASN:ND2	1.95	0.82
1:B:369:ARG:HE	1:B:370:ASN:HD21	1.24	0.82
1:B:51:ASN:HD21	1:B:53:SER:HB2	1.44	0.82
1:A:106:ASN:HD21	1:A:166:LYS:NZ	1.78	0.81
1:B:153:GLN:HB2	1:B:168:LEU:HD21	1.60	0.81
1:A:151:SER:H	1:A:169:ASN:HD21	1.25	0.81
1:A:158:VAL:HB	1:A:162:THR:HB	1.60	0.81
1:B:216:ARG:HB3	1:B:245:GLN:HG2	1.61	0.81
1:A:155:LEU:HB2	1:A:164:LYS:HG2	1.62	0.81
1:A:56:SER:N	1:A:59:TYR:HB2	1.97	0.80
1:B:309:ARG:HD3	1:B:309:ARG:N	1.93	0.80
1:B:104:PRO:HD2	1:B:171:GLY:O	1.82	0.79
1:B:481:SER:HB2	1:B:496:GLN:HB3	1.62	0.79
1:B:158:VAL:HB	1:B:162:THR:HB	1.64	0.79
1:A:254:LEU:O	1:A:292:LYS:HE3	1.83	0.79

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:56:SER:N	1:B:59:TYR:HB2	1.98	0.79
1:A:151:SER:N	1:A:169:ASN:HD21	1.80	0.79
1:B:90:ILE:HD12	1:B:130:TYR:HB2	1.64	0.79
1:B:52:PHE:O	1:B:86:LYS:HB2	1.82	0.78
1:A:369:ARG:HE	1:A:370:ASN:ND2	1.80	0.78
1:B:366:ASP:O	1:B:369:ARG:HG3	1.83	0.78
1:A:481:SER:HB2	1:A:496:GLN:HB3	1.65	0.78
1:B:51:ASN:ND2	1:B:53:SER:HB2	1.99	0.78
1:B:396:LYS:O	1:B:418:CYS:HB2	1.83	0.78
1:B:254:LEU:O	1:B:292:LYS:HE3	1.82	0.77
1:A:51:ASN:HD21	1:A:53:SER:HB2	1.49	0.77
1:A:90:ILE:HD12	1:A:130:TYR:HB2	1.65	0.77
1:B:139:ALA:HA	1:B:154:VAL:HG12	1.67	0.77
1:B:141:ARG:HH12	1:B:181:LEU:HB3	1.50	0.76
1:B:408:PRO:HG3	1:B:422:LEU:HD13	1.66	0.76
1:A:284:ILE:HG23	1:A:294:VAL:HG11	1.66	0.76
1:A:275:GLU:O	1:A:279:VAL:HG23	1.86	0.75
1:A:295:ILE:HG12	1:A:328:CYS:HB2	1.68	0.75
1:A:33:ASN:HD21	1:A:36:THR:H	1.32	0.75
1:B:496:GLN:HE21	1:B:498:SER:HB2	1.50	0.75
1:B:90:ILE:HB	1:B:179:VAL:HB	1.67	0.75
1:B:189:PRO:HB2	1:B:191:LEU:O	1.87	0.75
1:A:216:ARG:HB3	1:A:245:GLN:HG2	1.68	0.75
1:A:51:ASN:ND2	1:A:53:SER:HB2	2.02	0.74
1:A:52:PHE:O	1:A:86:LYS:HB2	1.87	0.74
1:A:125:ILE:HG12	1:A:126:MET:H	1.53	0.74
1:A:408:PRO:HG3	1:A:422:LEU:HD13	1.69	0.74
1:A:117:TYR:O	1:A:121:CYS:HB3	1.88	0.74
1:A:263:ALA:HA	2:A:1005:PGA:O2	1.87	0.74
1:B:369:ARG:HE	1:B:370:ASN:ND2	1.85	0.74
1:B:151:SER:H	1:B:169:ASN:HD21	1.36	0.74
1:A:366:ASP:O	1:A:369:ARG:HG3	1.88	0.74
1:A:496:GLN:HG3	1:A:497:VAL:N	2.02	0.73
1:A:346:THR:O	1:A:350:GLU:HG3	1.87	0.73
1:B:151:SER:N	1:B:169:ASN:HD21	1.86	0.73
1:B:117:TYR:O	1:B:121:CYS:HB3	1.89	0.73
1:B:155:LEU:HB2	1:B:164:LYS:HG2	1.71	0.73
1:B:496:GLN:HG3	1:B:497:VAL:H	1.53	0.73
1:B:372:THR:O	1:B:372:THR:HG22	1.87	0.73
1:B:496:GLN:HG3	1:B:497:VAL:N	2.04	0.72
1:B:398:ILE:HG23	1:B:420:ILE:HA	1.71	0.72
1:B:417:ASN:HD22	1:B:417:ASN:H	1.36	0.72

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:396:LYS:O	1:A:418:CYS:HB2	1.88	0.72
1:A:423:VAL:HG11	1:A:459:ARG:HB3	1.72	0.72
1:A:151:SER:H	1:A:169:ASN:ND2	1.88	0.72
1:B:295:ILE:HG12	1:B:328:CYS:HB2	1.70	0.72
1:B:125:ILE:HG12	1:B:126:MET:H	1.55	0.72
1:A:311:THR:OG1	1:A:314:GLU:HG3	1.90	0.71
1:A:398:ILE:HG23	1:A:420:ILE:HA	1.71	0.71
1:B:213:SER:HA	1:B:240:LYS:HD3	1.73	0.71
1:A:417:ASN:H	1:A:417:ASN:HD22	1.39	0.71
1:B:86:LYS:HG2	1:B:89:GLU:OE2	1.91	0.71
1:A:125:ILE:HG12	1:A:126:MET:N	2.06	0.71
1:A:496:GLN:HG3	1:A:497:VAL:H	1.55	0.71
1:A:372:THR:HG22	1:A:372:THR:O	1.88	0.71
1:B:346:THR:O	1:B:350:GLU:HG3	1.91	0.70
1:B:499:THR:HG22	1:B:500:VAL:N	2.07	0.70
1:B:79:LEU:HD12	1:B:80:ALA:H	1.55	0.70
1:A:263:ALA:C	2:A:1005:PGA:O2	2.29	0.70
1:A:153:GLN:HB2	1:A:168:LEU:HD21	1.74	0.70
1:B:456:VAL:O	1:B:460:ILE:HG13	1.92	0.70
1:A:496:GLN:HE21	1:A:498:SER:HB2	1.56	0.70
1:A:121:CYS:O	1:A:122:ASP:HB3	1.92	0.70
1:A:476:GLY:H	1:A:500:VAL:HB	1.57	0.70
1:B:67:ALA:HB1	1:B:79:LEU:HD21	1.74	0.70
1:B:284:ILE:HG23	1:B:294:VAL:HG11	1.73	0.70
1:A:48:VAL:HG23	1:A:79:LEU:HD11	1.72	0.69
1:A:106:ASN:HD21	1:A:166:LYS:HZ3	1.41	0.69
1:B:209:MET:O	1:B:210:VAL:HG23	1.91	0.69
1:B:243:ASN:HA	1:B:270:GLU:HG2	1.75	0.69
1:B:499:THR:HG22	1:B:500:VAL:H	1.56	0.69
1:A:104:PRO:HD2	1:A:171:GLY:O	1.93	0.69
1:A:86:LYS:HG2	1:A:89:GLU:OE2	1.93	0.68
1:A:141:ARG:HH12	1:A:181:LEU:HB3	1.58	0.68
1:A:101:TYR:HD2	1:A:172:LYS:HZ1	1.40	0.68
1:A:499:THR:HG22	1:A:500:VAL:H	1.59	0.68
1:A:369:ARG:NE	1:A:370:ASN:HD21	1.92	0.68
1:A:90:ILE:HG23	1:A:130:TYR:HB2	1.75	0.68
1:A:499:THR:HG22	1:A:500:VAL:N	2.08	0.68
1:A:247:VAL:C	1:A:249:ASN:H	1.97	0.68
1:B:125:ILE:HG12	1:B:126:MET:N	2.09	0.67
1:A:208:HIS:NE2	1:A:431:ARG:HD3	2.10	0.67
1:B:264:ARG:HH21	1:B:280:GLN:HE22	1.40	0.67
1:B:301:LEU:HD23	1:B:314:GLU:HB3	1.75	0.67

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:304:MET:HG3	1:A:310:PRO:HB3	1.76	0.67
1:A:213:SER:HA	1:A:240:LYS:HD3	1.77	0.67
1:B:124:LYS:N	1:B:124:LYS:HD2	2.10	0.67
1:B:311:THR:OG1	1:B:314:GLU:HG3	1.95	0.67
1:B:476:GLY:H	1:B:500:VAL:HB	1.59	0.67
1:A:243:ASN:HA	1:A:270:GLU:HG2	1.76	0.67
1:B:53:SER:HA	1:B:86:LYS:CG	2.24	0.66
1:B:475:LYS:HA	1:B:500:VAL:HB	1.78	0.66
1:B:360:ALA:HB1	1:B:363:PRO:HG2	1.78	0.66
1:B:275:GLU:O	1:B:279:VAL:HG23	1.95	0.66
1:B:133:ILE:HA	1:B:136:VAL:HG22	1.76	0.66
1:A:53:SER:HA	1:A:86:LYS:CG	2.24	0.66
1:A:84:ASP:CG	1:A:240:LYS:HZ2	1.98	0.66
1:A:33:ASN:ND2	1:A:36:THR:H	1.93	0.66
1:B:216:ARG:HG2	1:B:243:ASN:HD21	1.61	0.66
1:A:56:SER:H	1:A:59:TYR:HB2	1.60	0.65
1:B:53:SER:HA	1:B:86:LYS:HG3	1.77	0.65
1:A:216:ARG:HG2	1:A:243:ASN:HD21	1.61	0.65
1:A:483:GLN:HE21	1:A:483:GLN:N	1.94	0.65
1:A:53:SER:HA	1:A:86:LYS:HG3	1.77	0.65
1:A:3:ARG:HG2	1:A:7:LEU:HD22	1.77	0.65
1:B:139:ALA:HA	1:B:154:VAL:CG1	2.26	0.65
1:A:360:ALA:HB1	1:A:363:PRO:HG2	1.79	0.65
1:A:107:HIS:CE1	1:A:109:MET:HB3	2.32	0.65
1:A:390:VAL:CB	1:A:395:ALA:HB3	2.21	0.65
1:B:369:ARG:NE	1:B:370:ASN:HD21	1.94	0.65
1:B:56:SER:H	1:B:59:TYR:HB2	1.60	0.65
1:B:64:ILE:HG12	1:B:81:ILE:HG21	1.79	0.65
1:B:483:GLN:HE21	1:B:483:GLN:N	1.94	0.65
1:B:423:VAL:HG11	1:B:459:ARG:HB3	1.78	0.65
1:A:475:LYS:HA	1:A:500:VAL:HB	1.79	0.65
1:B:211:PHE:O	1:B:240:LYS:HD2	1.97	0.65
1:B:387:VAL:O	1:B:390:VAL:HG13	1.97	0.64
1:B:308:PRO:C	1:B:309:ARG:HD3	2.18	0.64
1:B:69:LYS:NZ	1:B:73:LEU:HG	2.12	0.64
1:B:33:ASN:HD21	1:B:36:THR:H	1.45	0.64
1:A:107:HIS:HE1	1:A:109:MET:HB3	1.61	0.64
1:B:112:THR:HG23	1:B:161:LYS:O	1.98	0.64
1:A:130:TYR:O	1:A:133:ILE:HG22	1.97	0.64
1:B:260:VAL:HG13	1:B:294:VAL:HG23	1.79	0.64
1:B:263:ALA:C	2:B:1006:PGA:O1	2.36	0.64
1:A:67:ALA:HB1	1:A:79:LEU:HD21	1.80	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:133:ILE:HA	1:A:136:VAL:HG22	1.79	0.63
1:A:196:LYS:HB3	1:A:200:ARG:HH21	1.63	0.63
1:B:247:VAL:C	1:B:249:ASN:H	2.02	0.63
1:B:208:HIS:NE2	1:B:431:ARG:HD3	2.13	0.63
1:A:375:PRO:O	1:A:376:THR:C	2.36	0.63
1:B:297:ALA:O	1:B:298:THR:HB	1.98	0.63
1:B:398:ILE:HD13	1:B:411:VAL:HG11	1.80	0.63
1:A:211:PHE:O	1:A:240:LYS:HD2	1.97	0.63
1:A:9:SER:O	1:A:11:ASN:N	2.31	0.63
1:B:3:ARG:HG2	1:B:7:LEU:HD22	1.78	0.63
1:B:304:MET:HG3	1:B:310:PRO:HB3	1.79	0.63
1:B:107:HIS:HE1	1:B:109:MET:HB3	1.63	0.63
1:A:64:ILE:HG12	1:A:81:ILE:HG21	1.80	0.63
1:A:257:THR:O	1:A:292:LYS:HD3	1.99	0.63
1:B:257:THR:O	1:B:292:LYS:HD3	1.98	0.63
1:A:429:ALA:HA	1:A:432:PHE:CZ	2.33	0.63
1:B:429:ALA:HA	1:B:432:PHE:CZ	2.34	0.63
1:A:209:MET:O	1:A:210:VAL:HG23	1.99	0.62
1:B:186:VAL:CG2	1:B:216:ARG:HE	2.06	0.62
1:A:113:THR:HG22	1:A:128:VAL:O	1.98	0.62
1:A:92:THR:HG22	1:A:177:LYS:H	1.64	0.62
1:A:141:ARG:HH22	1:A:181:LEU:HA	1.63	0.62
1:B:101:TYR:CD1	1:B:174:CYS:HA	2.33	0.62
1:B:92:THR:HG22	1:B:177:LYS:H	1.65	0.62
1:A:112:THR:HG23	1:A:161:LYS:O	1.99	0.62
1:A:110:ILE:HD11	1:A:162:THR:HG23	1.82	0.62
1:B:111:PHE:CZ	1:B:126:MET:SD	2.93	0.61
1:B:232:GLY:O	1:B:235:VAL:HG22	2.00	0.61
1:A:101:TYR:HB3	1:A:172:LYS:HG3	1.82	0.61
1:A:12:VAL:O	1:A:13:VAL:HB	2.00	0.61
1:A:180:ASN:O	1:A:182:PRO:HD3	2.01	0.61
1:B:113:THR:HG22	1:B:128:VAL:O	2.00	0.61
1:B:375:PRO:HB3	1:B:376:THR:OG1	1.99	0.61
1:B:56:SER:O	1:B:59:TYR:HB2	2.01	0.61
1:B:9:SER:O	1:B:11:ASN:N	2.33	0.61
1:B:151:SER:H	1:B:169:ASN:ND2	1.98	0.61
1:B:375:PRO:HB2	1:B:376:THR:HG1	1.66	0.60
1:B:141:ARG:HH22	1:B:181:LEU:HA	1.65	0.60
1:B:362:LEU:HB2	1:B:363:PRO:CD	2.31	0.60
1:B:2:SER:HB2	1:B:5:GLU:HB2	1.83	0.60
1:A:301:LEU:HD23	1:A:314:GLU:HB3	1.82	0.60
1:B:398:ILE:HG12	1:B:398:ILE:O	2.01	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:102:PRO:HB3	1:B:123:ASP:OD2	2.01	0.60
1:A:2:SER:HB2	1:A:5:GLU:HB2	1.82	0.60
1:A:93:GLY:HA3	1:A:127:TYR:HB3	1.82	0.60
1:A:297:ALA:O	1:A:298:THR:HB	2.02	0.60
1:A:112:THR:HB	1:A:125:ILE:HD11	1.82	0.60
1:B:107:HIS:CE1	1:B:109:MET:HB3	2.37	0.60
1:B:141:ARG:HH12	1:B:181:LEU:CB	2.14	0.60
1:A:403:THR:HG23	1:A:426:CYS:HB2	1.83	0.60
1:A:474:LYS:O	1:A:475:LYS:HB3	2.00	0.60
1:A:409:ARG:O	1:A:412:SER:HB3	2.01	0.60
1:B:396:LYS:O	1:B:419:PRO:HD2	2.02	0.60
1:B:475:LYS:HA	1:B:500:VAL:CG1	2.32	0.59
1:B:242:GLU:HA	1:B:267:LEU:HD13	1.84	0.59
1:B:48:VAL:HG23	1:B:79:LEU:HD11	1.83	0.59
1:B:151:SER:HB2	1:B:168:LEU:HB2	1.84	0.59
1:A:65:ASP:HA	1:A:68:ARG:HB2	1.84	0.59
1:B:95:THR:HA	1:B:121:CYS:O	2.03	0.59
1:A:41:ARG:NH2	1:A:74:TYR:O	2.36	0.59
1:A:56:SER:O	1:A:59:TYR:HB2	2.03	0.59
1:A:242:GLU:HA	1:A:267:LEU:HD13	1.84	0.59
1:A:300:MET:O	1:A:314:GLU:HB3	2.03	0.59
1:B:15:GLY:H	1:B:357:GLN:HE22	1.51	0.59
1:A:139:ALA:HA	1:A:154:VAL:HG12	1.85	0.58
1:A:398:ILE:HD13	1:A:411:VAL:HG11	1.83	0.58
1:A:387:VAL:O	1:A:390:VAL:HG13	2.03	0.58
1:A:15:GLY:H	1:A:357:GLN:HE22	1.52	0.58
1:B:156:GLU:HB2	1:B:164:LYS:HE2	1.86	0.58
1:A:151:SER:HB2	1:A:168:LEU:HB2	1.85	0.58
1:B:121:CYS:O	1:B:122:ASP:HB3	2.03	0.58
1:B:474:LYS:O	1:B:475:LYS:HB3	2.04	0.58
1:B:477:ASP:O	1:B:478:THR:HB	2.03	0.58
1:B:195:ASP:HA	1:B:198:ASP:OD2	2.03	0.58
1:B:151:SER:HB3	1:B:168:LEU:HD12	1.85	0.58
1:A:12:VAL:HG23	1:A:13:VAL:HG23	1.85	0.58
1:B:94:THR:HA	1:B:175:SER:HB2	1.85	0.58
1:A:101:TYR:CE1	1:A:174:CYS:HB3	2.39	0.58
1:A:69:LYS:NZ	1:A:73:LEU:HG	2.18	0.58
1:B:105:PRO:HD2	1:B:107:HIS:CD2	2.39	0.57
1:A:94:THR:HA	1:A:175:SER:HB2	1.84	0.57
1:B:417:ASN:N	1:B:417:ASN:HD22	2.01	0.57
1:A:303:SER:C	1:A:305:THR:H	2.07	0.57
1:B:69:LYS:HZ2	1:B:73:LEU:HG	1.68	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:91:ARG:HG2	1:A:91:ARG:HH11	1.69	0.57
1:A:475:LYS:HA	1:A:500:VAL:CG1	2.34	0.57
1:B:480:VAL:HG13	1:B:496:GLN:O	2.03	0.57
1:B:110:ILE:HD11	1:B:162:THR:HG23	1.86	0.57
1:B:264:ARG:HB2	1:B:297:ALA:O	2.05	0.57
1:A:106:ASN:ND2	1:A:166:LYS:NZ	2.50	0.57
1:A:112:THR:HA	1:A:161:LYS:O	2.05	0.57
1:A:189:PRO:HB2	1:A:191:LEU:O	2.05	0.57
1:A:101:TYR:CD1	1:A:174:CYS:HA	2.39	0.57
1:B:403:THR:HG23	1:B:426:CYS:HB2	1.87	0.57
1:B:112:THR:HB	1:B:125:ILE:HD11	1.87	0.57
1:B:65:ASP:HA	1:B:68:ARG:HB2	1.87	0.57
1:B:431:ARG:O	1:B:434:HIS:HD2	1.88	0.57
1:A:456:VAL:O	1:A:460:ILE:HG13	2.05	0.57
1:A:141:ARG:HH12	1:A:181:LEU:CB	2.16	0.56
1:B:33:ASN:ND2	1:B:36:THR:H	2.02	0.56
1:B:409:ARG:O	1:B:412:SER:HB3	2.05	0.56
1:A:92:THR:O	1:A:176:HIS:HD2	1.87	0.56
1:A:141:ARG:HE	1:A:142:ILE:N	2.04	0.56
1:A:84:ASP:OD1	1:A:240:LYS:NZ	2.37	0.56
1:B:180:ASN:O	1:B:182:PRO:HD3	2.05	0.56
1:B:110:ILE:HD13	1:B:125:ILE:HG13	1.87	0.56
1:A:401:LEU:N	1:A:401:LEU:HD12	2.21	0.56
1:A:264:ARG:HB2	1:A:297:ALA:O	2.06	0.55
1:A:110:ILE:HD13	1:A:125:ILE:HG13	1.88	0.55
1:A:101:TYR:HD2	1:A:172:LYS:NZ	2.03	0.55
1:A:247:VAL:O	1:A:249:ASN:N	2.39	0.55
1:A:153:GLN:HG3	1:A:168:LEU:HD21	1.88	0.55
1:A:396:LYS:O	1:A:419:PRO:HD2	2.06	0.55
1:A:79:LEU:HD12	1:A:80:ALA:H	1.71	0.55
1:A:232:GLY:O	1:A:235:VAL:HG22	2.05	0.55
1:B:362:LEU:HB2	1:B:363:PRO:HD3	1.87	0.55
1:B:112:THR:HA	1:B:161:LYS:O	2.07	0.55
1:A:480:VAL:HG13	1:A:496:GLN:O	2.05	0.55
1:B:67:ALA:HB1	1:B:79:LEU:CD2	2.36	0.55
1:B:141:ARG:HE	1:B:142:ILE:N	2.05	0.55
1:B:300:MET:O	1:B:314:GLU:HB3	2.06	0.55
1:B:153:GLN:CB	1:B:168:LEU:HD21	2.32	0.55
1:A:67:ALA:HB1	1:A:79:LEU:CD2	2.36	0.55
1:A:195:ASP:HA	1:A:198:ASP:OD2	2.07	0.55
1:A:444:PHE:CE2	1:A:459:ARG:HG2	2.42	0.55
1:B:79:LEU:HD12	1:B:80:ALA:N	2.22	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:12:VAL:O	1:B:13:VAL:HB	2.06	0.55
1:A:114:ASP:O	1:A:115:ASP:HB2	2.06	0.54
1:A:477:ASP:O	1:A:478:THR:HB	2.07	0.54
1:B:102:PRO:O	1:B:173:ILE:HG22	2.06	0.54
1:A:398:ILE:O	1:A:398:ILE:HG12	2.07	0.54
1:A:208:HIS:CE1	1:A:431:ARG:HD3	2.43	0.54
1:A:401:LEU:HD23	1:A:459:ARG:HD2	1.89	0.54
1:B:460:ILE:O	1:B:464:ILE:HG13	2.08	0.54
1:A:57:TYR:O	1:A:61:LYS:HB2	2.08	0.54
1:B:399:ILE:HA	1:B:421:ILE:O	2.07	0.54
1:B:495:LEU:C	1:B:495:LEU:HD23	2.28	0.54
1:B:86:LYS:NZ	1:B:89:GLU:HG3	2.22	0.54
1:B:41:ARG:NH2	1:B:74:TYR:O	2.40	0.54
1:B:141:ARG:HH21	1:B:143:ILE:HA	1.73	0.54
1:A:476:GLY:N	1:A:500:VAL:HB	2.21	0.54
1:B:208:HIS:CE1	1:B:431:ARG:HD3	2.43	0.54
1:B:101:TYR:HD2	1:B:172:LYS:HZ1	1.56	0.54
1:B:110:ILE:HD11	1:B:112:THR:OG1	2.07	0.54
1:B:45:LEU:HD21	1:B:48:VAL:CG2	2.38	0.54
1:A:47:ILE:HG12	1:A:80:ALA:HB3	1.89	0.54
1:B:402:SER:HB3	1:B:422:LEU:HD11	1.89	0.54
1:A:362:LEU:HB2	1:A:363:PRO:CD	2.38	0.53
1:B:159:ASP:HB2	1:B:162:THR:OG1	2.09	0.53
1:A:52:PHE:CD2	1:A:198:ASP:HB3	2.44	0.53
1:A:146:ASP:OD1	1:A:177:LYS:HD2	2.08	0.53
1:B:444:PHE:CE2	1:B:459:ARG:HG2	2.43	0.53
1:A:431:ARG:O	1:A:434:HIS:HD2	1.90	0.53
1:B:130:TYR:O	1:B:133:ILE:HG22	2.08	0.53
1:B:47:ILE:HG12	1:B:80:ALA:HB3	1.91	0.53
1:A:399:ILE:HA	1:A:421:ILE:O	2.09	0.53
1:B:45:LEU:HD21	1:B:48:VAL:HG22	1.91	0.53
1:A:427:PRO:O	1:A:431:ARG:HG3	2.09	0.53
1:B:273:ALA:N	1:B:274:PRO:HD2	2.24	0.53
1:A:453:THR:HG22	1:A:453:THR:O	2.09	0.53
1:A:478:THR:O	1:A:478:THR:HG23	2.09	0.53
1:B:445:GLU:O	1:B:446:LYS:CB	2.56	0.53
1:B:196:LYS:HB3	1:B:200:ARG:HH21	1.74	0.53
1:A:69:LYS:HZ2	1:A:73:LEU:HG	1.73	0.53
1:B:114:ASP:O	1:B:115:ASP:HB2	2.08	0.53
1:A:112:THR:CB	1:A:125:ILE:HD11	2.39	0.53
1:A:52:PHE:HZ	1:A:201:PHE:CE2	2.27	0.53
1:B:153:GLN:HG3	1:B:168:LEU:HD21	1.91	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:496:GLN:CG	1:A:497:VAL:H	2.22	0.52
1:B:303:SER:C	1:B:305:THR:H	2.11	0.52
1:B:362:LEU:N	1:B:362:LEU:HD13	2.24	0.52
1:B:86:LYS:HZ2	1:B:89:GLU:HG3	1.75	0.52
1:B:476:GLY:N	1:B:500:VAL:HB	2.23	0.52
1:A:45:LEU:HD21	1:A:48:VAL:HG22	1.91	0.52
1:A:102:PRO:HB3	1:A:123:ASP:OD2	2.09	0.52
1:B:427:PRO:O	1:B:431:ARG:HG3	2.09	0.52
1:B:10:LEU:O	1:B:11:ASN:C	2.47	0.52
1:B:84:ASP:CG	1:B:240:LYS:HZ2	2.12	0.52
1:B:496:GLN:CG	1:B:497:VAL:H	2.19	0.52
1:A:247:VAL:C	1:A:249:ASN:N	2.63	0.52
1:A:1:MET:O	1:A:3:ARG:N	2.43	0.52
1:B:105:PRO:CD	1:B:107:HIS:HD2	2.23	0.52
1:B:123:ASP:HB3	1:B:124:LYS:HD2	1.92	0.52
1:B:425:ARG:HA	1:B:444:PHE:O	2.10	0.52
1:B:398:ILE:HG22	1:B:419:PRO:O	2.10	0.52
1:A:103:ILE:HA	1:A:172:LYS:HA	1.92	0.52
1:B:6:ARG:NE	1:B:360:ALA:HB2	2.25	0.52
1:B:95:THR:H	1:B:175:SER:HB3	1.74	0.52
1:B:456:VAL:HG12	1:B:459:ARG:NH2	2.25	0.52
1:A:6:ARG:NE	1:A:360:ALA:HB2	2.24	0.52
1:A:151:SER:HB3	1:A:168:LEU:HD12	1.91	0.52
1:B:1:MET:O	1:B:3:ARG:N	2.43	0.52
1:B:429:ALA:HA	1:B:432:PHE:CE2	2.45	0.52
1:A:111:PHE:CD1	1:A:165:VAL:HG21	2.45	0.52
1:A:264:ARG:HE	1:A:280:GLN:NE2	2.08	0.51
1:A:425:ARG:HA	1:A:444:PHE:O	2.09	0.51
1:A:45:LEU:HD21	1:A:48:VAL:CG2	2.41	0.51
1:A:362:LEU:HB2	1:A:363:PRO:HD3	1.91	0.51
1:A:105:PRO:HD2	1:A:107:HIS:CD2	2.45	0.51
1:A:95:THR:HA	1:A:121:CYS:O	2.09	0.51
1:A:159:ASP:HB2	1:A:162:THR:OG1	2.10	0.51
1:B:496:GLN:CG	1:B:497:VAL:N	2.73	0.51
1:A:223:THR:O	1:A:227:VAL:HG23	2.10	0.51
1:B:410:LEU:O	1:B:413:LYS:HB3	2.10	0.51
1:A:52:PHE:HZ	1:A:201:PHE:HE2	1.59	0.51
1:B:213:SER:CA	1:B:240:LYS:HD3	2.41	0.51
1:A:86:LYS:HZ2	1:A:89:GLU:HG3	1.74	0.51
1:B:122:ASP:O	1:B:124:LYS:N	2.42	0.51
1:A:361:TYR:CG	1:A:417:ASN:HB3	2.45	0.51
1:A:496:GLN:CG	1:A:497:VAL:N	2.73	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:223:THR:O	1:B:227:VAL:HG23	2.11	0.51
1:B:28:GLY:N	1:B:336:ALA:HA	2.26	0.51
1:B:106:ASN:HD21	1:B:166:LYS:NZ	2.09	0.51
1:A:495:LEU:C	1:A:495:LEU:HD23	2.31	0.51
1:A:446:LYS:O	1:A:447:GLU:CB	2.58	0.50
1:A:19:ARG:NH1	1:A:21:THR:O	2.42	0.50
1:B:382:VAL:HG21	1:B:493:ASN:ND2	2.26	0.50
1:A:429:ALA:HA	1:A:432:PHE:CE2	2.45	0.50
1:B:401:LEU:HD12	1:B:401:LEU:N	2.26	0.50
1:B:279:VAL:HG12	1:B:279:VAL:O	2.10	0.50
1:A:432:PHE:O	1:A:435:LEU:HB2	2.12	0.50
1:B:432:PHE:C	1:B:432:PHE:CD1	2.84	0.50
1:A:156:GLU:HB2	1:A:164:LYS:HE2	1.94	0.50
1:A:247:VAL:HG13	1:A:248:ASN:N	2.26	0.50
1:B:52:PHE:CD2	1:B:198:ASP:HB3	2.47	0.50
1:A:445:GLU:O	1:A:446:LYS:CB	2.60	0.50
1:B:475:LYS:HA	1:B:500:VAL:CB	2.40	0.50
1:B:359:ILE:HG22	1:B:360:ALA:N	2.26	0.50
1:B:23:ILE:HG21	1:B:345:VAL:HG13	1.94	0.50
1:B:398:ILE:CG2	1:B:420:ILE:HA	2.39	0.50
1:A:398:ILE:CG2	1:A:420:ILE:HA	2.40	0.50
1:A:410:LEU:O	1:A:413:LYS:HB3	2.12	0.50
1:A:34:PRO:O	1:A:38:VAL:HG23	2.11	0.50
1:B:374:LYS:N	1:B:375:PRO:CB	2.67	0.49
1:A:399:ILE:O	1:A:481:SER:HA	2.11	0.49
1:B:361:TYR:CG	1:B:417:ASN:HB3	2.46	0.49
1:B:390:VAL:CB	1:B:395:ALA:HB3	2.29	0.49
1:B:141:ARG:NH2	1:B:143:ILE:HA	2.28	0.49
1:B:439:VAL:O	1:B:441:PRO:HD3	2.12	0.49
1:B:483:GLN:HG2	1:B:484:GLY:N	2.25	0.49
1:A:425:ARG:HH21	1:A:447:GLU:N	2.09	0.49
1:B:90:ILE:CB	1:B:179:VAL:HB	2.40	0.49
1:B:57:TYR:O	1:B:61:LYS:HB2	2.12	0.49
1:A:273:ALA:N	1:A:274:PRO:HD2	2.26	0.49
1:A:279:VAL:O	1:A:283:LEU:HD22	2.12	0.49
1:B:425:ARG:HH21	1:B:447:GLU:N	2.10	0.49
1:B:81:ILE:HD13	1:B:81:ILE:N	2.28	0.49
1:A:457:GLU:H	1:A:457:GLU:CD	2.16	0.49
1:B:93:GLY:HA3	1:B:127:TYR:HB3	1.95	0.49
1:B:188:LEU:HD23	1:B:189:PRO:HD2	1.94	0.49
1:B:112:THR:CB	1:B:125:ILE:HD11	2.42	0.49
1:B:499:THR:CG2	1:B:500:VAL:N	2.75	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:141:ARG:HH21	1:A:143:ILE:HA	1.76	0.49
1:B:266:ASP:O	1:B:269:ILE:N	2.41	0.49
1:B:85:THR:HG22	1:B:87:GLY:N	2.27	0.49
1:A:31:THR:O	1:A:37:LEU:HD22	2.13	0.49
1:B:103:ILE:HG12	1:B:172:LYS:HB2	1.94	0.49
1:A:386:ALA:O	1:A:389:ALA:HB3	2.12	0.49
1:A:28:GLY:N	1:A:336:ALA:HA	2.27	0.49
1:A:122:ASP:O	1:A:124:LYS:N	2.43	0.49
1:B:264:ARG:HE	1:B:280:GLN:NE2	2.11	0.49
1:B:146:ASP:O	1:B:149:VAL:HG13	2.12	0.49
1:B:499:THR:CG2	1:B:500:VAL:H	2.25	0.49
1:B:105:PRO:HD2	1:B:107:HIS:HD2	1.78	0.49
1:A:460:ILE:O	1:A:464:ILE:HG13	2.13	0.49
1:A:212:ALA:O	1:A:240:LYS:HB2	2.12	0.49
1:A:86:LYS:NZ	1:A:89:GLU:HG3	2.27	0.48
1:A:89:GLU:O	1:A:90:ILE:HD13	2.12	0.48
1:A:266:ASP:O	1:A:269:ILE:N	2.45	0.48
1:A:141:ARG:HH12	1:A:181:LEU:CA	2.26	0.48
1:B:247:VAL:HG13	1:B:248:ASN:N	2.28	0.48
1:A:305:THR:HG23	1:A:338:GLY:HA2	1.95	0.48
1:A:231:GLN:NE2	1:A:231:GLN:H	2.11	0.48
1:B:258:ASP:O	1:B:293:PRO:HD2	2.13	0.48
1:B:400:VAL:O	1:B:422:LEU:HA	2.14	0.48
1:A:139:ALA:HA	1:A:154:VAL:CG1	2.43	0.48
1:B:457:GLU:H	1:B:457:GLU:CD	2.17	0.48
1:B:375:PRO:CB	1:B:376:THR:HG1	2.23	0.48
1:B:52:PHE:HZ	1:B:201:PHE:CE2	2.32	0.48
1:A:301:LEU:O	1:A:304:MET:HB2	2.13	0.48
1:A:475:LYS:HA	1:A:500:VAL:CB	2.42	0.48
1:B:113:THR:CG2	1:B:128:VAL:HG23	2.43	0.48
1:B:422:LEU:HD23	1:B:441:PRO:HB3	1.94	0.48
1:B:12:VAL:HG23	1:B:13:VAL:HG23	1.94	0.48
1:B:114:ASP:OD2	1:B:116:LYS:HG3	2.12	0.48
1:B:186:VAL:HG23	1:B:216:ARG:NE	2.05	0.48
1:B:305:THR:HG23	1:B:338:GLY:HA2	1.96	0.48
1:B:386:ALA:O	1:B:389:ALA:HB3	2.14	0.48
1:B:477:ASP:O	1:B:478:THR:CB	2.61	0.48
1:B:342:ILE:CD1	1:B:342:ILE:H	2.10	0.48
1:A:94:THR:HG22	1:A:175:SER:HB2	1.95	0.48
1:B:401:LEU:HD23	1:B:459:ARG:HD2	1.94	0.48
1:A:432:PHE:CD1	1:A:432:PHE:C	2.87	0.48
1:B:478:THR:HG23	1:B:478:THR:O	2.13	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:141:ARG:HH12	1:B:181:LEU:CA	2.27	0.48
1:A:175:SER:O	1:A:177:LYS:HG2	2.14	0.48
1:B:100:ASP:OD1	1:B:122:ASP:OD2	2.31	0.48
1:A:279:VAL:O	1:A:279:VAL:HG12	2.13	0.47
1:A:146:ASP:O	1:A:149:VAL:HG22	2.14	0.47
1:B:399:ILE:O	1:B:481:SER:HA	2.15	0.47
1:B:84:ASP:OD1	1:B:240:LYS:NZ	2.42	0.47
1:B:86:LYS:HZ3	1:B:89:GLU:CD	2.17	0.47
1:B:110:ILE:HG13	1:B:162:THR:HG22	1.95	0.47
1:B:188:LEU:CD2	1:B:189:PRO:HD2	2.44	0.47
1:A:402:SER:HB3	1:A:422:LEU:HD11	1.96	0.47
1:A:95:THR:H	1:A:175:SER:HB3	1.79	0.47
1:B:314:GLU:O	1:B:318:VAL:HG23	2.14	0.47
1:B:215:ILE:HG13	1:B:239:VAL:HG13	1.95	0.47
1:A:425:ARG:NH2	1:A:447:GLU:N	2.62	0.47
1:A:361:TYR:HB3	1:A:391:PHE:CZ	2.49	0.47
1:A:456:VAL:HG12	1:A:459:ARG:NH2	2.29	0.47
1:B:247:VAL:C	1:B:249:ASN:N	2.67	0.47
1:B:247:VAL:O	1:B:249:ASN:N	2.45	0.47
1:A:153:GLN:CB	1:A:168:LEU:HD21	2.43	0.47
1:B:141:ARG:HE	1:B:142:ILE:H	1.63	0.47
1:B:156:GLU:OE1	1:B:164:LYS:NZ	2.48	0.47
1:A:10:LEU:O	1:A:11:ASN:C	2.52	0.47
1:B:231:GLN:NE2	1:B:231:GLN:H	2.13	0.47
1:B:415:ARG:NH1	1:B:437:ARG:HB3	2.29	0.47
1:A:404:SER:HB2	1:A:406:THR:OG1	2.15	0.47
1:B:130:TYR:CE1	1:B:132:ASN:HB2	2.49	0.47
1:A:297:ALA:HB1	1:A:330:MET:HE1	1.96	0.47
1:A:48:VAL:HG23	1:A:79:LEU:CD1	2.42	0.47
1:B:33:ASN:O	1:B:37:LEU:HB2	2.15	0.47
1:B:19:ARG:NH1	1:B:21:THR:O	2.41	0.47
1:B:175:SER:O	1:B:177:LYS:HG2	2.15	0.47
1:B:247:VAL:CG1	1:B:248:ASN:N	2.78	0.47
1:A:91:ARG:HG2	1:A:91:ARG:NH1	2.29	0.47
1:B:467:ALA:HB1	1:B:473:LEU:HD13	1.96	0.47
1:B:425:ARG:NH2	1:B:447:GLU:N	2.63	0.46
1:A:400:VAL:O	1:A:422:LEU:HA	2.15	0.46
1:A:114:ASP:C	1:A:116:LYS:H	2.18	0.46
1:B:453:THR:HG22	1:B:453:THR:O	2.15	0.46
1:A:106:ASN:ND2	1:A:106:ASN:O	2.48	0.46
1:B:92:THR:O	1:B:176:HIS:HD2	1.99	0.46
1:A:361:TYR:HB3	1:A:391:PHE:HZ	1.80	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:110:ILE:HD13	1:A:110:ILE:O	2.15	0.46
1:A:415:ARG:NH1	1:A:437:ARG:HB3	2.31	0.46
1:B:41:ARG:HA	1:B:45:LEU:HB3	1.98	0.46
1:A:382:VAL:HG21	1:A:493:ASN:ND2	2.31	0.46
1:B:88:PRO:HB2	1:B:188:LEU:HD13	1.96	0.46
1:B:111:PHE:CE2	1:B:126:MET:SD	3.09	0.46
1:A:88:PRO:HG3	1:A:189:PRO:O	2.16	0.46
1:A:90:ILE:HB	1:A:179:VAL:HB	1.98	0.46
1:A:456:VAL:HG12	1:A:459:ARG:CZ	2.46	0.46
1:A:401:LEU:HD21	1:A:459:ARG:HB2	1.97	0.46
1:A:362:LEU:H	1:A:362:LEU:HD22	1.81	0.46
1:A:115:ASP:HA	1:A:118:ALA:HB2	1.97	0.46
1:B:329:VAL:HB	1:B:348:MET:SD	2.55	0.46
1:A:216:ARG:HH11	1:A:216:ARG:HG3	1.80	0.46
1:B:112:THR:HG22	1:B:113:THR:N	2.31	0.46
1:A:90:ILE:HD12	1:A:130:TYR:CB	2.39	0.46
1:A:499:THR:CG2	1:A:500:VAL:N	2.77	0.46
1:B:87:GLY:O	1:B:89:GLU:HG2	2.15	0.45
1:B:110:ILE:CD1	1:B:112:THR:OG1	2.64	0.45
1:A:192:SER:H	1:A:195:ASP:HB2	1.81	0.45
1:B:446:LYS:O	1:B:447:GLU:CB	2.63	0.45
1:A:359:ILE:HG22	1:A:360:ALA:N	2.32	0.45
1:A:126:MET:HE1	1:A:173:ILE:HG13	1.98	0.45
1:B:185:ASP:OD1	1:B:245:GLN:NE2	2.49	0.45
1:B:133:ILE:O	1:B:137:ILE:HB	2.17	0.45
1:B:156:GLU:HB2	1:B:164:LYS:CE	2.46	0.45
1:B:327:ASP:HA	1:B:437:ARG:HB2	1.98	0.45
1:B:243:ASN:HA	1:B:270:GLU:CG	2.43	0.45
1:B:484:GLY:HA2	1:B:493:ASN:HA	1.99	0.45
1:A:110:ILE:HD11	1:A:112:THR:OG1	2.16	0.45
1:A:141:ARG:NH2	1:A:143:ILE:HA	2.32	0.45
1:A:221:VAL:O	1:A:224:ILE:HG23	2.16	0.45
1:B:304:MET:HE1	1:B:343:ASN:CB	2.46	0.45
1:A:85:THR:HG22	1:A:87:GLY:N	2.31	0.45
1:A:216:ARG:HD3	1:A:245:GLN:NE2	2.31	0.45
1:A:247:VAL:CG1	1:A:248:ASN:N	2.79	0.45
1:A:439:VAL:O	1:A:441:PRO:HD3	2.16	0.45
1:B:102:PRO:CG	1:B:122:ASP:OD2	2.65	0.45
1:A:213:SER:CA	1:A:240:LYS:HD3	2.44	0.45
1:A:362:LEU:N	1:A:362:LEU:HD13	2.32	0.45
1:A:50:MET:HE3	1:A:63:VAL:HB	1.98	0.45
1:B:90:ILE:CD1	1:B:130:TYR:HB2	2.41	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:86:LYS:HZ3	1:A:89:GLU:CD	2.19	0.45
1:B:88:PRO:HB2	1:B:188:LEU:HB3	1.99	0.45
1:A:243:ASN:HA	1:A:270:GLU:CG	2.45	0.45
1:A:141:ARG:HE	1:A:142:ILE:H	1.64	0.45
1:B:103:ILE:HA	1:B:172:LYS:HA	1.98	0.45
1:A:113:THR:CG2	1:A:128:VAL:HG23	2.46	0.45
1:B:50:MET:HE3	1:B:63:VAL:HB	1.98	0.45
1:B:105:PRO:O	1:B:106:ASN:HB2	2.17	0.44
1:B:158:VAL:CB	1:B:162:THR:HB	2.43	0.44
1:A:260:VAL:HG13	1:A:294:VAL:HG23	1.98	0.44
1:B:153:GLN:CG	1:B:168:LEU:HD21	2.47	0.44
1:A:159:ASP:HB3	1:A:160:ASP:H	1.60	0.44
1:A:444:PHE:HB2	1:A:462:PHE:CD2	2.53	0.44
1:B:45:LEU:HD11	1:B:47:ILE:O	2.18	0.44
1:A:99:VAL:HG12	1:A:101:TYR:CZ	2.53	0.44
1:B:330:MET:HG3	1:B:331:LEU:H	1.83	0.44
1:A:81:ILE:HD13	1:A:81:ILE:N	2.32	0.44
1:B:28:GLY:HA3	1:B:336:ALA:O	2.17	0.44
1:A:204:LYS:HE3	1:A:204:LYS:HB2	1.78	0.44
1:A:56:SER:H	1:A:59:TYR:CB	2.27	0.44
1:B:158:VAL:O	1:B:162:THR:HB	2.18	0.44
1:A:87:GLY:O	1:A:89:GLU:HG2	2.17	0.44
1:B:48:VAL:HG23	1:B:79:LEU:CD1	2.47	0.44
1:A:141:ARG:NH2	1:A:181:LEU:HA	2.32	0.44
1:A:99:VAL:O	1:A:99:VAL:HG12	2.17	0.44
1:A:397:ALA:HB2	1:A:473:LEU:HD11	1.99	0.44
1:A:477:ASP:O	1:A:478:THR:CB	2.66	0.44
1:A:216:ARG:NH1	1:A:216:ARG:HG3	2.32	0.44
1:A:124:LYS:N	1:A:124:LYS:HD2	2.33	0.44
1:B:41:ARG:HD3	1:B:79:LEU:HD22	1.99	0.44
1:A:103:ILE:HG12	1:A:172:LYS:HB2	2.00	0.44
1:B:415:ARG:HH11	1:B:437:ARG:HB3	1.82	0.44
1:A:89:GLU:HG2	1:A:89:GLU:H	1.62	0.44
1:B:94:THR:HG22	1:B:175:SER:HB2	2.00	0.44
1:B:456:VAL:HG12	1:B:459:ARG:CZ	2.48	0.44
1:B:145:VAL:O	1:B:149:VAL:HG22	2.17	0.44
1:B:216:ARG:HD3	1:B:245:GLN:NE2	2.33	0.44
1:B:407:THR:N	1:B:408:PRO:HD2	2.32	0.44
1:A:483:GLN:HG2	1:A:484:GLY:N	2.33	0.44
1:B:301:LEU:O	1:B:304:MET:HB2	2.18	0.44
1:B:398:ILE:CD1	1:B:411:VAL:HG11	2.47	0.44
1:B:91:ARG:HH11	1:B:91:ARG:HG2	1.83	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:40:LEU:HD12	1:A:40:LEU:HA	1.81	0.44
1:A:263:ALA:O	1:A:267:LEU:HB2	2.18	0.43
1:A:133:ILE:O	1:A:137:ILE:HB	2.17	0.43
1:B:444:PHE:HB2	1:B:462:PHE:CD2	2.53	0.43
1:A:242:GLU:HA	1:A:267:LEU:HB2	1.99	0.43
1:B:191:LEU:HD22	1:B:195:ASP:HB3	1.98	0.43
1:A:158:VAL:O	1:A:162:THR:HB	2.18	0.43
1:B:62:SER:OG	1:B:63:VAL:N	2.51	0.43
1:A:112:THR:HG22	1:A:113:THR:N	2.33	0.43
1:B:90:ILE:O	1:B:178:GLY:HA2	2.18	0.43
1:A:208:HIS:HA	1:A:428:ARG:HH12	1.84	0.43
1:B:115:ASP:HA	1:B:118:ALA:HB2	2.00	0.43
1:A:467:ALA:HB1	1:A:473:LEU:HD13	2.00	0.43
1:B:304:MET:HG3	1:B:310:PRO:CB	2.48	0.43
1:B:122:ASP:OD1	1:B:122:ASP:C	2.56	0.43
1:B:156:GLU:HB2	1:B:164:LYS:NZ	2.34	0.43
1:A:398:ILE:HG22	1:A:419:PRO:O	2.18	0.43
1:A:423:VAL:HG21	1:A:459:ARG:O	2.19	0.43
1:A:425:ARG:HG2	1:A:444:PHE:O	2.19	0.43
1:A:141:ARG:HH22	1:A:181:LEU:HD23	1.84	0.43
1:B:37:LEU:HA	1:B:37:LEU:HD13	1.77	0.43
1:B:242:GLU:HA	1:B:267:LEU:HB2	2.01	0.43
1:A:331:LEU:HA	1:A:331:LEU:HD23	1.90	0.43
1:B:279:VAL:O	1:B:283:LEU:HD22	2.19	0.43
1:A:102:PRO:O	1:A:173:ILE:HG22	2.19	0.43
1:A:493:ASN:N	1:A:493:ASN:OD1	2.51	0.43
1:A:180:ASN:ND2	1:A:270:GLU:OE1	2.52	0.43
1:B:401:LEU:HD21	1:B:459:ARG:HB2	2.00	0.43
1:B:65:ASP:O	1:B:69:LYS:HB2	2.18	0.43
1:B:243:ASN:CA	1:B:270:GLU:HG2	2.44	0.43
1:A:484:GLY:HA2	1:A:493:ASN:HA	2.01	0.43
1:B:222:LEU:O	1:B:225:ARG:HB3	2.18	0.43
1:B:88:PRO:HG3	1:B:189:PRO:O	2.19	0.43
1:B:496:GLN:HE21	1:B:498:SER:CB	2.24	0.43
1:B:113:THR:HG22	1:B:128:VAL:HG23	2.00	0.43
1:B:133:ILE:O	1:B:137:ILE:N	2.47	0.43
1:A:90:ILE:HG23	1:A:130:TYR:CB	2.48	0.43
1:B:391:PHE:HA	1:B:391:PHE:HD1	1.73	0.43
1:B:146:ASP:O	1:B:149:VAL:HG22	2.19	0.43
1:A:327:ASP:HA	1:A:437:ARG:HB2	2.00	0.43
1:A:155:LEU:CB	1:A:164:LYS:HG2	2.39	0.43
1:B:181:LEU:HB3	1:B:184:THR:HB	2.00	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:117:TYR:CD1	1:B:117:TYR:N	2.86	0.43
1:A:60:HIS:C	1:A:62:SER:N	2.72	0.43
1:B:387:VAL:HG21	1:B:414:TYR:HB2	2.01	0.42
1:A:281:LYS:HZ3	1:A:324:ASP:CG	2.22	0.42
1:A:30:LYS:C	1:A:30:LYS:HD3	2.40	0.42
1:B:30:LYS:HD3	1:B:30:LYS:C	2.38	0.42
1:A:153:GLN:CG	1:A:168:LEU:HD21	2.49	0.42
1:B:496:GLN:HG3	1:B:498:SER:H	1.84	0.42
1:A:41:ARG:HA	1:A:45:LEU:HB3	2.01	0.42
1:A:45:LEU:HD11	1:A:47:ILE:O	2.19	0.42
1:B:115:ASP:OD1	1:B:131:LYS:HE3	2.19	0.42
1:B:8:THR:HG22	1:B:8:THR:O	2.19	0.42
1:A:65:ASP:O	1:A:69:LYS:HB2	2.19	0.42
1:B:466:LYS:HZ2	1:B:469:GLU:CD	2.21	0.42
1:B:109:MET:HG3	1:B:110:ILE:N	2.34	0.42
1:B:56:SER:H	1:B:59:TYR:CB	2.28	0.42
1:A:133:ILE:HD11	1:A:137:ILE:HD12	2.01	0.42
1:A:88:PRO:HD3	1:A:190:ALA:O	2.19	0.42
1:A:417:ASN:N	1:A:417:ASN:HD22	2.04	0.42
1:A:314:GLU:O	1:A:318:VAL:HG23	2.18	0.42
1:B:432:PHE:C	1:B:432:PHE:HD1	2.22	0.42
1:B:194:LYS:O	1:B:197:GLU:HB2	2.20	0.42
1:A:276:VAL:O	1:A:280:GLN:N	2.52	0.42
1:B:56:SER:OG	1:B:59:TYR:HD1	2.02	0.42
1:B:7:LEU:HA	1:B:10:LEU:HD22	2.01	0.42
1:B:361:TYR:HB3	1:B:391:PHE:CZ	2.54	0.42
1:A:433:SER:C	1:A:435:LEU:H	2.23	0.42
1:A:110:ILE:HG13	1:A:162:THR:HG22	2.01	0.42
1:B:407:THR:N	1:B:408:PRO:CD	2.82	0.42
1:A:499:THR:CG2	1:A:500:VAL:H	2.27	0.42
1:A:143:ILE:HB	1:A:152:PHE:HB2	2.02	0.42
1:A:244:GLN:HA	1:A:247:VAL:HG12	2.01	0.42
1:A:342:ILE:H	1:A:342:ILE:CD1	2.12	0.42
1:B:89:GLU:H	1:B:89:GLU:HG2	1.62	0.42
1:A:37:LEU:HD13	1:A:37:LEU:HA	1.79	0.42
1:A:361:TYR:CD1	1:A:417:ASN:HB3	2.55	0.42
1:A:41:ARG:HG2	1:A:45:LEU:HD23	2.02	0.42
1:B:432:PHE:O	1:B:435:LEU:HB2	2.20	0.42
1:B:410:LEU:HA	1:B:410:LEU:HD12	1.90	0.42
1:A:23:ILE:HG21	1:A:345:VAL:HG13	2.02	0.42
1:B:497:VAL:HG22	1:B:497:VAL:O	2.20	0.42
1:B:289:LEU:HD12	1:B:289:LEU:HA	1.90	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:106:ASN:HD21	1:A:166:LYS:HZ1	1.60	0.41
1:A:243:ASN:CA	1:A:270:GLU:HG2	2.46	0.41
1:A:110:ILE:HD11	1:A:162:THR:CG2	2.50	0.41
1:A:113:THR:HB	1:A:131:LYS:NZ	2.35	0.41
1:B:132:ASN:O	1:B:133:ILE:C	2.58	0.41
1:A:481:SER:CB	1:A:496:GLN:HB3	2.42	0.41
1:A:180:ASN:OD1	1:A:269:ILE:HG21	2.20	0.41
1:A:304:MET:HG3	1:A:310:PRO:CB	2.48	0.41
1:A:407:THR:N	1:A:408:PRO:HD2	2.35	0.41
1:A:387:VAL:HG21	1:A:414:TYR:HB2	2.02	0.41
1:B:153:GLN:HB2	1:B:168:LEU:CD2	2.41	0.41
1:A:254:LEU:HD13	1:A:290:ALA:HB2	2.02	0.41
1:A:199:LEU:O	1:A:203:VAL:HG23	2.21	0.41
1:A:33:ASN:O	1:A:37:LEU:HB2	2.19	0.41
1:B:67:ALA:HA	1:B:70:SER:OG	2.20	0.41
1:B:263:ALA:O	1:B:267:LEU:HB2	2.21	0.41
1:A:12:VAL:O	1:A:13:VAL:CB	2.68	0.41
1:A:194:LYS:O	1:A:197:GLU:HB2	2.20	0.41
1:B:88:PRO:CB	1:B:188:LEU:HB3	2.50	0.41
1:A:130:TYR:CE1	1:A:132:ASN:HB2	2.56	0.41
1:A:122:ASP:C	1:A:122:ASP:OD1	2.58	0.41
1:A:24:ILE:HG12	1:A:47:ILE:HB	2.02	0.41
1:B:297:ALA:HB1	1:B:330:MET:HE1	2.03	0.41
1:A:472:ILE:HG22	1:A:473:LEU:HD12	2.02	0.41
1:B:416:PRO:CG	1:B:420:ILE:HD11	2.51	0.41
1:A:66:ASN:HA	1:A:66:ASN:HD22	1.64	0.41
1:B:144:TYR:HA	1:B:150:LEU:O	2.21	0.41
1:B:239:VAL:HG23	1:B:257:THR:OG1	2.21	0.41
1:A:100:ASP:OD1	1:A:122:ASP:OD2	2.38	0.41
1:A:46:ASN:O	1:A:79:LEU:HD12	2.21	0.41
1:A:132:ASN:O	1:A:133:ILE:C	2.59	0.41
1:A:86:LYS:HD3	1:A:86:LYS:O	2.21	0.41
1:B:330:MET:CG	1:B:331:LEU:H	2.34	0.41
1:B:52:PHE:HZ	1:B:201:PHE:HE2	1.67	0.41
1:A:195:ASP:O	1:A:199:LEU:HG	2.20	0.41
1:A:180:ASN:C	1:A:182:PRO:HD3	2.42	0.41
1:B:41:ARG:HG2	1:B:45:LEU:HD23	2.02	0.41
1:B:330:MET:CG	1:B:331:LEU:N	2.84	0.41
1:B:362:LEU:HD22	1:B:362:LEU:H	1.86	0.41
1:B:362:LEU:H	1:B:362:LEU:HD13	1.85	0.41
1:B:208:HIS:HA	1:B:428:ARG:HH12	1.86	0.41
1:B:144:TYR:HB3	1:B:148:GLY:HA2	2.02	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:281:LYS:HG3	1:B:281:LYS:H	1.55	0.41
1:A:215:ILE:HG13	1:A:239:VAL:HG13	2.03	0.41
1:B:342:ILE:HG12	1:B:343:ASN:N	2.36	0.41
1:B:69:LYS:HA	1:B:72:GLU:OE2	2.20	0.41
1:A:234:ASP:N	1:A:234:ASP:OD1	2.53	0.41
1:B:243:ASN:CB	1:B:270:GLU:HG2	2.51	0.40
1:A:276:VAL:O	1:A:277:LEU:C	2.59	0.40
1:A:125:ILE:CG1	1:A:126:MET:N	2.82	0.40
1:B:262:VAL:O	1:B:264:ARG:N	2.53	0.40
1:B:493:ASN:N	1:B:493:ASN:OD1	2.54	0.40
1:B:47:ILE:HG22	1:B:48:VAL:N	2.36	0.40
1:A:303:SER:C	1:A:305:THR:N	2.73	0.40
1:A:415:ARG:HH11	1:A:437:ARG:HB3	1.86	0.40
1:A:289:LEU:HD11	1:A:371:CYS:CB	2.52	0.40
1:A:26:THR:O	1:A:335:THR:HB	2.20	0.40
1:A:496:GLN:HG3	1:A:498:SER:H	1.86	0.40
1:B:361:TYR:CD1	1:B:417:ASN:HB3	2.56	0.40
1:B:263:ALA:O	2:B:1006:PGA:O1	2.39	0.40
1:B:1:MET:HG2	1:B:2:SER:N	2.36	0.40
1:A:28:GLY:HA3	1:A:336:ALA:O	2.21	0.40
1:B:151:SER:CB	1:B:168:LEU:HB2	2.48	0.40
1:A:372:THR:CG2	1:A:372:THR:O	2.60	0.40
1:A:222:LEU:O	1:A:225:ARG:HB3	2.22	0.40
1:B:220:ASP:O	1:B:224:ILE:CG2	2.56	0.40
1:A:191:LEU:HD22	1:A:195:ASP:HB3	2.04	0.40
1:A:201:PHE:O	1:A:205:ASN:ND2	2.52	0.40
1:B:102:PRO:HG2	1:B:122:ASP:OD2	2.22	0.40

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

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:309:ARG:NH2	1:B:268:GLY:O[1_545]	2.06	0.14

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	481/500 (96%)	382 (79%)	65 (14%)	34 (7%)	2	9
1	B	481/500 (96%)	377 (78%)	71 (15%)	33 (7%)	2	9
All	All	962/1000 (96%)	759 (79%)	136 (14%)	67 (7%)	2	9

All (67) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	10	LEU
1	A	12	VAL
1	A	96	THR
1	A	99	VAL
1	A	106	ASN
1	A	375	PRO
1	A	398	ILE
1	A	446	LYS
1	A	478	THR
1	A	497	VAL
1	B	10	LEU
1	B	96	THR
1	B	99	VAL
1	B	106	ASN
1	B	374	LYS
1	B	375	PRO
1	B	398	ILE
1	B	446	LYS
1	B	478	THR
1	B	497	VAL
1	A	2	SER
1	A	85	THR
1	A	87	GLY
1	A	123	ASP
1	A	146	ASP
1	A	176	HIS
1	A	265	GLY
1	A	374	LYS
1	B	2	SER
1	B	12	VAL
1	B	85	THR
1	B	123	ASP
1	B	146	ASP
1	B	176	HIS
1	B	265	GLY

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Mol	Chain	Res	Type
1	A	11	ASN
1	A	248	ASN
1	A	249	ASN
1	A	494	THR
1	B	11	ASN
1	B	87	GLY
1	B	248	ASN
1	B	249	ASN
1	B	494	THR
1	B	495	LEU
1	A	13	VAL
1	A	105	PRO
1	A	122	ASP
1	A	170	ALA
1	A	298	THR
1	A	304	MET
1	B	13	VAL
1	B	105	PRO
1	B	118	ALA
1	B	255	LYS
1	B	298	THR
1	A	255	LYS
1	A	475	LYS
1	A	495	LEU
1	B	170	ALA
1	B	304	MET
1	B	475	LYS
1	A	110	ILE
1	A	210	VAL
1	A	372	THR
1	B	372	THR
1	B	210	VAL

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	412/423 (97%)	337 (82%)	75 (18%)	2 13

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	B	412/423 (97%)	339 (82%)	73 (18%)	3 14
All	All	824/846 (97%)	676 (82%)	148 (18%)	2 13

All (148) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	2	SER
1	A	7	LEU
1	A	10	LEU
1	A	22	SER
1	A	30	LYS
1	A	31	THR
1	A	33	ASN
1	A	37	LEU
1	A	40	LEU
1	A	50	MET
1	A	58	GLU
1	A	61	LYS
1	A	65	ASP
1	A	66	ASN
1	A	68	ARG
1	A	69	LYS
1	A	70	SER
1	A	77	ARG
1	A	83	LEU
1	A	89	GLU
1	A	91	ARG
1	A	94	THR
1	A	95	THR
1	A	96	THR
1	A	98	ASP
1	A	99	VAL
1	A	100	ASP
1	A	110	ILE
1	A	122	ASP
1	A	128	VAL
1	A	138	SER
1	A	141	ARG
1	A	150	LEU
1	A	153	GLN
1	A	156	GLU

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Mol	Chain	Res	Type
1	A	160	ASP
1	A	162	THR
1	A	169	ASN
1	A	173	ILE
1	A	186	VAL
1	A	224	ILE
1	A	225	ARG
1	A	230	GLU
1	A	245	GLN
1	A	254	LEU
1	A	262	VAL
1	A	281	LYS
1	A	283	LEU
1	A	309	ARG
1	A	311	THR
1	A	316	SER
1	A	323	LEU
1	A	329	VAL
1	A	337	LYS
1	A	342	ILE
1	A	362	LEU
1	A	369	ARG
1	A	376	THR
1	A	390	VAL
1	A	391	PHE
1	A	400	VAL
1	A	404	SER
1	A	407	THR
1	A	410	LEU
1	A	415	ARG
1	A	417	ASN
1	A	431	ARG
1	A	432	PHE
1	A	433	SER
1	A	445	GLU
1	A	455	ASP
1	A	474	LYS
1	A	483	GLN
1	A	498	SER
1	B	1	MET
1	B	2	SER
1	B	7	LEU

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Mol	Chain	Res	Type
1	B	10	LEU
1	B	22	SER
1	B	30	LYS
1	B	31	THR
1	B	33	ASN
1	B	37	LEU
1	B	40	LEU
1	B	50	MET
1	B	58	GLU
1	B	61	LYS
1	B	65	ASP
1	B	66	ASN
1	B	68	ARG
1	B	69	LYS
1	B	70	SER
1	B	83	LEU
1	B	89	GLU
1	B	94	THR
1	B	95	THR
1	B	96	THR
1	B	98	ASP
1	B	100	ASP
1	B	110	ILE
1	B	122	ASP
1	B	126	MET
1	B	128	VAL
1	B	138	SER
1	B	141	ARG
1	B	150	LEU
1	B	153	GLN
1	B	156	GLU
1	B	160	ASP
1	B	162	THR
1	B	169	ASN
1	B	173	ILE
1	B	186	VAL
1	B	224	ILE
1	B	225	ARG
1	B	245	GLN
1	B	254	LEU
1	B	281	LYS
1	B	283	LEU

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Mol	Chain	Res	Type
1	B	309	ARG
1	B	311	THR
1	B	316	SER
1	B	323	LEU
1	B	324	ASP
1	B	329	VAL
1	B	337	LYS
1	B	342	ILE
1	B	362	LEU
1	B	369	ARG
1	B	376	THR
1	B	390	VAL
1	B	391	PHE
1	B	400	VAL
1	B	404	SER
1	B	407	THR
1	B	410	LEU
1	B	415	ARG
1	B	417	ASN
1	B	428	ARG
1	B	431	ARG
1	B	432	PHE
1	B	433	SER
1	B	445	GLU
1	B	455	ASP
1	B	474	LYS
1	B	483	GLN
1	B	498	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	ASN
1	A	60	HIS
1	A	66	ASN
1	A	106	ASN
1	A	169	ASN
1	A	176	HIS
1	A	231	GLN
1	A	243	ASN
1	A	244	GLN
1	A	245	GLN

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Mol	Chain	Res	Type
1	A	280	GLN
1	A	343	ASN
1	A	357	GLN
1	A	370	ASN
1	A	417	ASN
1	A	434	HIS
1	A	483	GLN
1	A	493	ASN
1	A	496	GLN
1	B	33	ASN
1	B	60	HIS
1	B	66	ASN
1	B	106	ASN
1	B	132	ASN
1	B	169	ASN
1	B	176	HIS
1	B	231	GLN
1	B	243	ASN
1	B	244	GLN
1	B	280	GLN
1	B	343	ASN
1	B	357	GLN
1	B	370	ASN
1	B	417	ASN
1	B	434	HIS
1	B	483	GLN
1	B	493	ASN
1	B	496	GLN

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

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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$
2	PGA	A	1005	1,3,4	8,8,8	3.78	2 (25%)	11,11,11	2.78	5 (45%)
2	PGA	B	1006	3,4	8,8,8	2.55	2 (25%)	11,11,11	2.84	5 (45%)

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
2	PGA	A	1005	1,3,4	-	0/6/6/6	0/0/0/0
2	PGA	B	1006	3,4	-	0/6/6/6	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1005	PGA	O1P-C2	-9.24	1.36	1.43
2	B	1006	PGA	O1P-C2	-4.89	1.39	1.43
2	A	1005	PGA	P-O3P	4.35	1.70	1.54
2	B	1006	PGA	P-O3P	4.21	1.70	1.54

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1006	PGA	O1P-P-O2P	6.57	125.98	106.71
2	A	1005	PGA	O1P-P-O2P	6.27	125.08	106.71
2	B	1006	PGA	O1P-C2-C1	3.53	116.24	110.54
2	A	1005	PGA	O1P-C2-C1	3.42	116.06	110.54
2	B	1006	PGA	P-O1P-C2	3.40	133.49	121.22
2	A	1005	PGA	P-O1P-C2	2.99	132.00	121.22
2	A	1005	PGA	O3P-P-O1P	2.80	114.39	106.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1005	PGA	O2-C1-O1	-2.41	117.17	123.30
2	B	1006	PGA	O3P-P-O1P	2.29	112.97	106.65
2	B	1006	PGA	O2-C1-O1	-2.24	117.60	123.30

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.