



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 09:24 am GMT

PDB ID : 4BY0
Title : Crystal structure of Trypanosoma cruzi CYP51 bound to the inhibitor (R)-N-(3-(1H-indol-3-yl)-1-oxo-1-(pyridin-4-ylamino)propan-2-yl)-3,3'-difluoro-(1,1'-biphenyl)-4-carboxamide
Authors : Choi, J.Y.; Calvet, C.M.; Vierira, D.F.; Gunatilleke, S.S.; Cameron, M.D.; McKerrow, J.H.; Podust, L.M.; Roush, W.R.
Deposited on : 2013-07-16
Resolution : 3.10 Å(reported)

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

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

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

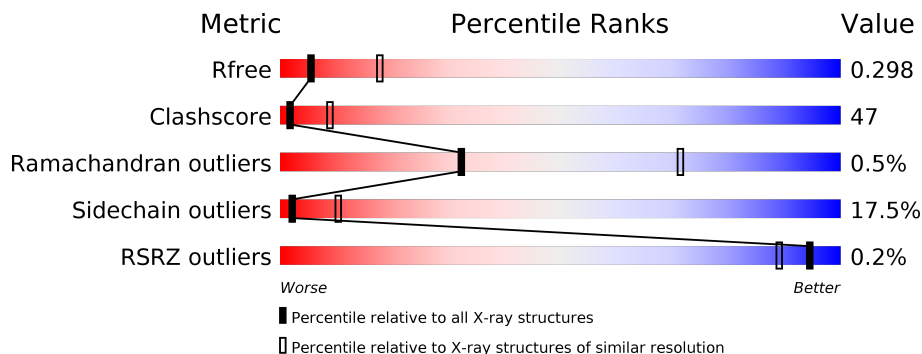
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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}	100719	1001 (3.12-3.08)
Clashscore	112137	1099 (3.12-3.08)
Ramachandran outliers	110173	1057 (3.12-3.08)
Sidechain outliers	110143	1057 (3.12-3.08)
RSRZ outliers	101464	1006 (3.12-3.08)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	467	
1	B	467	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	5PS	A	1460	-	-	X	-
3	5PS	B	1460	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6834 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called STEROL 14-ALPHA DEMETHYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	441	Total	C	N	O	S	0	0	0
			3389	2162	589	611	27			
1	B	442	Total	C	N	O	S	0	0	0
			3272	2077	570	598	27			

There are 34 discrepancies between the modelled and reference sequences:

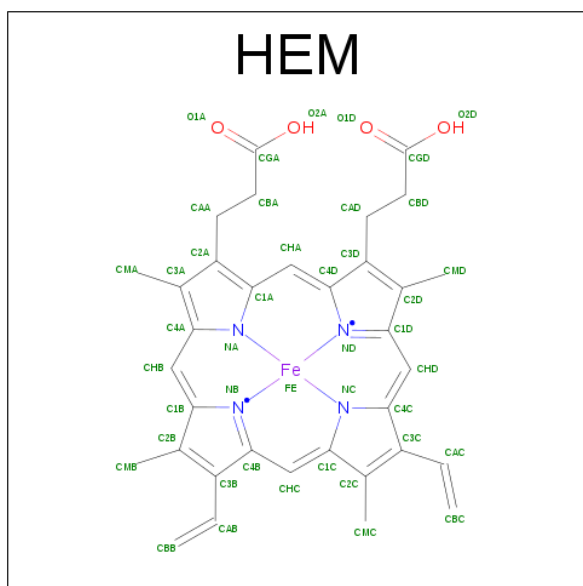
Chain	Residue	Modelled	Actual	Comment	Reference
A	21	MET	-	EXPRESSION TAG	UNP Q5I4E1
A	22	ALA	-	EXPRESSION TAG	UNP Q5I4E1
A	23	LYS	-	EXPRESSION TAG	UNP Q5I4E1
A	24	LYS	-	EXPRESSION TAG	UNP Q5I4E1
A	25	THR	-	EXPRESSION TAG	UNP Q5I4E1
A	26	SER	-	EXPRESSION TAG	UNP Q5I4E1
A	27	SER	-	EXPRESSION TAG	UNP Q5I4E1
A	28	LYS	-	EXPRESSION TAG	UNP Q5I4E1
A	29	GLY	-	EXPRESSION TAG	UNP Q5I4E1
A	30	LYS	-	EXPRESSION TAG	UNP Q5I4E1
A	31	LEU	-	EXPRESSION TAG	UNP Q5I4E1
A	482	HIS	-	EXPRESSION TAG	UNP Q5I4E1
A	483	HIS	-	EXPRESSION TAG	UNP Q5I4E1
A	484	HIS	-	EXPRESSION TAG	UNP Q5I4E1
A	485	HIS	-	EXPRESSION TAG	UNP Q5I4E1
A	486	HIS	-	EXPRESSION TAG	UNP Q5I4E1
A	487	HIS	-	EXPRESSION TAG	UNP Q5I4E1
B	21	MET	-	EXPRESSION TAG	UNP Q5I4E1
B	22	ALA	-	EXPRESSION TAG	UNP Q5I4E1
B	23	LYS	-	EXPRESSION TAG	UNP Q5I4E1
B	24	LYS	-	EXPRESSION TAG	UNP Q5I4E1
B	25	THR	-	EXPRESSION TAG	UNP Q5I4E1
B	26	SER	-	EXPRESSION TAG	UNP Q5I4E1
B	27	SER	-	EXPRESSION TAG	UNP Q5I4E1
B	28	LYS	-	EXPRESSION TAG	UNP Q5I4E1

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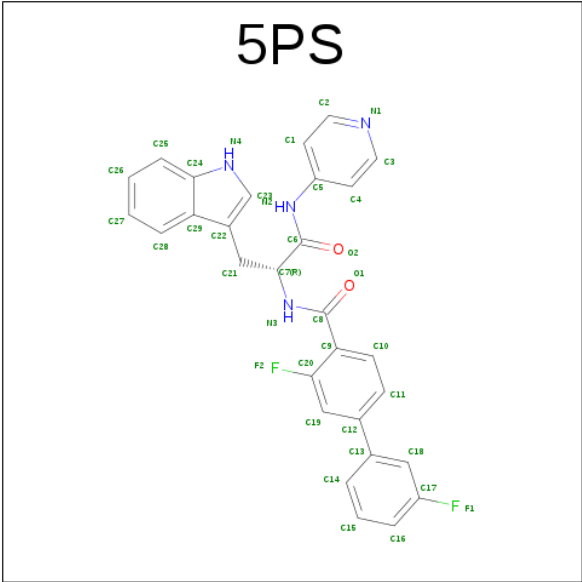
Chain	Residue	Modelled	Actual	Comment	Reference
B	29	GLY	-	EXPRESSION TAG	UNP Q5I4E1
B	30	LYS	-	EXPRESSION TAG	UNP Q5I4E1
B	31	LEU	-	EXPRESSION TAG	UNP Q5I4E1
B	482	HIS	-	EXPRESSION TAG	UNP Q5I4E1
B	483	HIS	-	EXPRESSION TAG	UNP Q5I4E1
B	484	HIS	-	EXPRESSION TAG	UNP Q5I4E1
B	485	HIS	-	EXPRESSION TAG	UNP Q5I4E1
B	486	HIS	-	EXPRESSION TAG	UNP Q5I4E1
B	487	HIS	-	EXPRESSION TAG	UNP Q5I4E1

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 3 is (R)-N-(3-(1H-INDOL-3-YL)-1-OXO-1-(PYRIDIN-4-YLAMINO)PROPAN-2-YL)-3,3'-DIFLUORO-(1,1'-BIPHENYL)-4-CARBOXAMIDE (three-letter code: 5PS) (formula: C₂₉H₂₂F₂N₄O₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			37	29	2	4	2		
3	B	1	Total	C	F	N	O	0	0
			37	29	2	4	2		

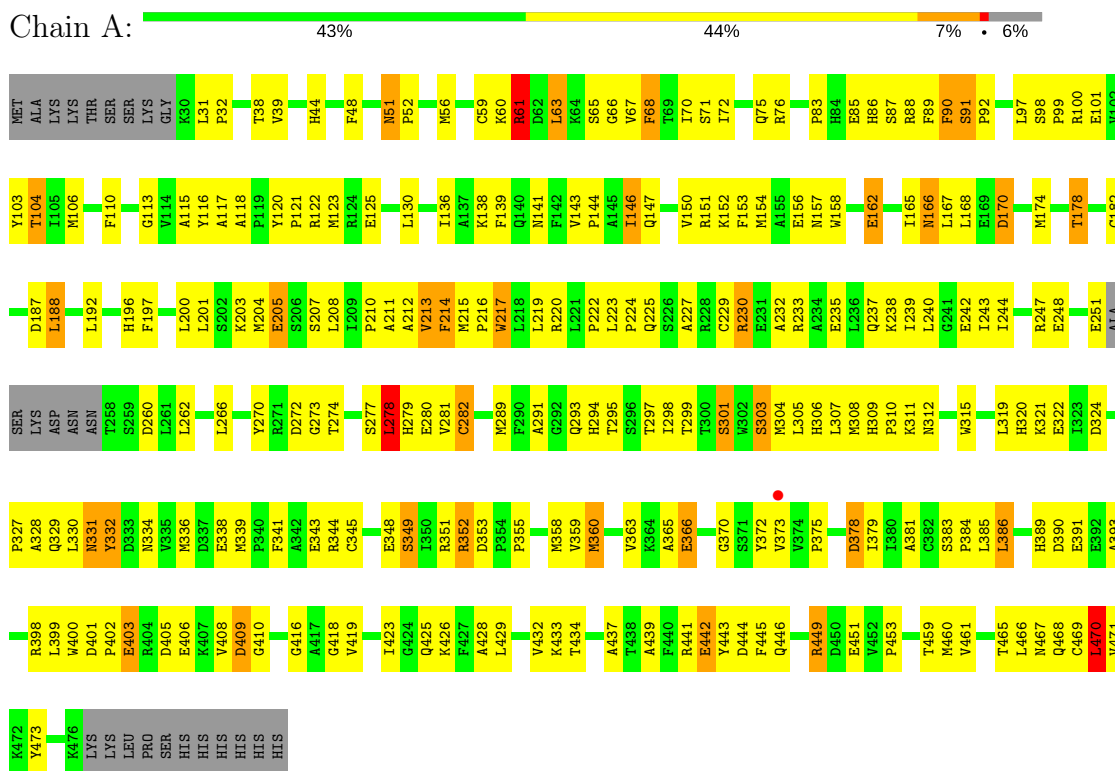
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	8	Total	O	0	0
			8	8		
4	B	5	Total	O	0	0
			5	5		

3 Residue-property plots

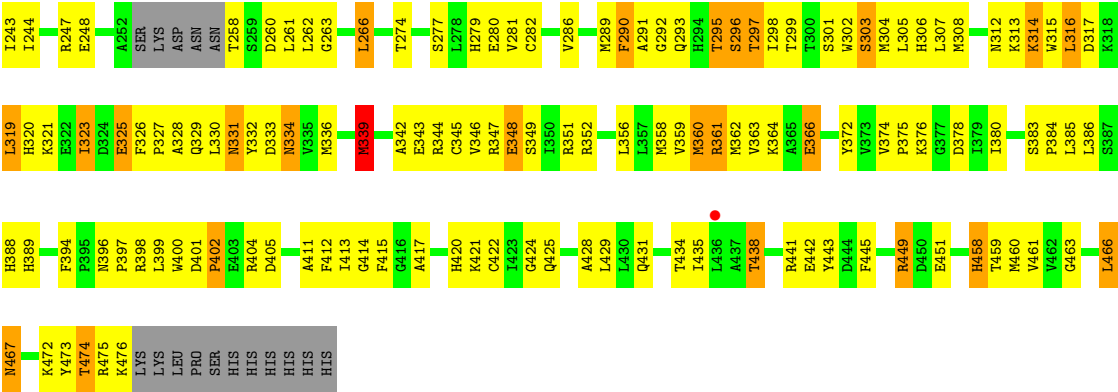
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

• Molecule 1: STEROL 14-ALPHA DEMETHYLASE



• Molecule 1: STEROL 14-ALPHA DEMETHYLASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	124.17Å 124.17Å 119.84Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	80.16 – 3.10 80.04 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (80.16-3.10) 99.9 (80.04-3.10)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 3.13Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.232 , 0.297 0.228 , 0.298	Depositor DCC
R_{free} test set	1013 reflections (5.40%)	DCC
Wilson B-factor (Å ²)	85.9	Xtriage
Anisotropy	0.146	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 83.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6834	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, 5PS

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	0.61	0/3469	0.78	4/4710 (0.1%)
1	B	0.55	0/3350	0.73	0/4562
All	All	0.58	0/6819	0.75	4/9272 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	61	ARG	NE-CZ-NH1	8.11	124.36	120.30
1	A	222	PRO	N-CA-CB	6.54	111.14	103.30
1	A	470	LEU	CA-CB-CG	-5.24	103.25	115.30
1	A	449	ARG	N-CA-C	-5.10	97.23	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3389	0	3249	255	0
1	B	3272	0	3007	350	0
2	A	43	0	30	13	0
2	B	43	0	30	16	0
3	A	37	0	22	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	37	0	22	5	0
4	A	8	0	0	0	0
4	B	5	0	0	1	0
All	All	6834	0	6360	620	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 47.

All (620) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:35:TYR:CD1	1:B:36:PRO:HD2	1.53	1.42
1:A:59:CYS:O	1:A:63:LEU:CD1	1.78	1.29
1:B:326:PHE:O	1:B:441:ARG:NH2	1.66	1.26
1:B:110:PHE:O	1:B:233:ARG:NH2	1.63	1.26
1:A:363:VAL:HG23	1:A:378:ASP:O	1.33	1.25
1:A:366:GLU:OE2	1:A:373:VAL:HG11	1.23	1.25
1:A:449:ARG:NE	1:A:468:GLN:OE1	1.71	1.22
1:B:344:ARG:NH2	1:B:405:ASP:OD1	1.71	1.21
1:B:66:GLY:O	1:B:81:GLY:HA3	1.11	1.21
1:A:366:GLU:OE2	1:A:373:VAL:CG1	1.88	1.20
1:A:308:MET:CE	1:A:445:PHE:HB3	1.72	1.20
1:B:172:GLY:CA	1:B:297:THR:HG21	1.69	1.19
1:B:66:GLY:O	1:B:81:GLY:CA	1.91	1.19
1:B:89:PHE:HD2	1:B:90:PHE:CE1	1.62	1.16
1:B:89:PHE:HD2	1:B:90:PHE:CD1	1.62	1.16
1:A:110:PHE:O	1:A:233:ARG:NH2	1.78	1.16
1:B:89:PHE:CD2	1:B:90:PHE:CE1	2.34	1.15
1:A:147:GLN:OE1	1:A:330:LEU:HB2	1.43	1.15
1:B:33:PRO:HB2	1:B:63:LEU:HD22	1.30	1.12
1:A:308:MET:CE	1:A:445:PHE:CB	2.28	1.11
1:B:83:PRO:O	1:B:86:HIS:HB2	1.48	1.11
1:B:344:ARG:NH1	1:B:405:ASP:OD1	1.84	1.09
1:B:344:ARG:CZ	1:B:405:ASP:OD1	2.01	1.09
1:B:89:PHE:CD2	1:B:90:PHE:CD1	2.41	1.08
1:A:59:CYS:C	1:A:63:LEU:HD12	1.72	1.08
1:A:308:MET:HE1	1:A:445:PHE:HB3	1.36	1.07
1:B:111:GLY:O	1:B:114:VAL:HG23	1.54	1.07
1:B:172:GLY:HA2	1:B:297:THR:CG2	1.84	1.06
1:B:188:LEU:HD13	1:B:243:ILE:HD12	1.35	1.06
1:B:331:ASN:ND2	1:B:334:ASN:OD1	1.90	1.05

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:277:SER:O	1:B:281:VAL:HG23	1.57	1.04
1:B:172:GLY:CA	1:B:297:THR:CG2	2.35	1.04
1:A:59:CYS:O	1:A:63:LEU:HD12	0.84	1.01
1:B:35:TYR:CD1	1:B:36:PRO:CD	2.43	1.01
1:B:344:ARG:HH22	1:B:405:ASP:CG	1.64	1.01
1:B:172:GLY:HA2	1:B:297:THR:HG21	1.40	1.00
1:B:215:MET:O	1:B:219:LEU:N	1.94	1.00
1:A:211:ALA:O	1:A:215:MET:N	1.94	1.00
1:B:420:HIS:ND1	2:B:1450:HEM:O1D	1.96	0.99
1:A:224:PRO:HB2	1:A:227:ALA:HB2	1.44	0.99
1:B:442:GLU:O	1:B:476:LYS:HB2	1.60	0.98
1:B:32:PRO:HB3	1:B:372:TYR:CD1	1.99	0.97
1:B:328:ALA:HB3	1:B:329:GLN:NE2	1.78	0.97
1:A:360:MET:N	1:A:360:MET:SD	2.38	0.96
1:B:360:MET:O	1:B:361:ARG:NH1	1.97	0.95
1:B:102:VAL:HG23	1:B:103:TYR:CD1	2.01	0.95
1:A:235:GLU:O	1:A:239:ILE:HG13	1.65	0.95
1:A:363:VAL:CG2	1:A:378:ASP:O	2.16	0.94
1:A:214:PHE:HB2	1:A:217:TRP:CZ3	2.02	0.93
1:B:361:ARG:HH22	2:B:1450:HEM:CGA	1.81	0.93
1:B:313:LYS:HA	1:B:316:LEU:HB3	1.50	0.93
1:B:172:GLY:N	1:B:297:THR:CG2	2.32	0.92
1:B:32:PRO:HA	1:B:372:TYR:CE1	2.04	0.92
1:B:33:PRO:CB	1:B:63:LEU:HD22	1.98	0.92
1:B:299:THR:O	1:B:303:SER:OG	1.87	0.92
1:B:147:GLN:O	1:B:151:ARG:N	2.03	0.91
1:A:308:MET:HE2	1:A:445:PHE:CB	1.98	0.91
1:B:139:PHE:HD2	1:B:332:TYR:HE2	1.16	0.91
1:A:48:PHE:O	1:A:52:PRO:HD3	1.69	0.91
1:A:449:ARG:CZ	1:A:468:GLN:OE1	2.20	0.89
1:A:211:ALA:C	1:A:215:MET:CB	2.40	0.89
1:B:277:SER:OG	1:B:280:GLU:HG3	1.72	0.89
1:B:175:ILE:CD1	1:B:296:SER:HB2	2.03	0.89
1:B:32:PRO:HA	1:B:372:TYR:HE1	1.38	0.89
1:B:361:ARG:NH2	2:B:1450:HEM:O1A	2.05	0.89
1:A:213:VAL:O	1:A:379:ILE:HD12	1.73	0.88
1:A:449:ARG:NH1	1:A:451:GLU:O	2.07	0.88
1:A:308:MET:CE	1:A:445:PHE:HB2	2.04	0.87
1:B:247:ARG:NH2	1:B:260:ASP:OD2	2.07	0.87
2:A:1450:HEM:C1A	3:A:1460:5PS:H2	2.10	0.86
1:A:339:MET:HE1	1:A:437:ALA:HB2	1.56	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:313:LYS:O	1:B:317:ASP:N	2.07	0.86
1:A:348:GLU:OE2	1:A:351:ARG:NH2	2.09	0.85
1:B:323:ILE:HA	1:B:326:PHE:CE2	2.10	0.85
1:B:86:HIS:CE1	1:B:413:ILE:HD13	2.10	0.85
1:A:465:THR:HB	1:A:468:GLN:HG3	1.57	0.85
1:A:211:ALA:HB1	1:A:215:MET:HA	1.59	0.85
1:A:143:VAL:HG21	1:A:332:TYR:HA	1.58	0.84
1:B:188:LEU:CD1	1:B:243:ILE:HD12	2.06	0.84
1:B:146:ILE:O	1:B:150:VAL:HG23	1.77	0.84
1:A:98:SER:OG	1:A:101:GLU:OE1	1.95	0.84
1:A:113:GLY:O	1:A:118:ALA:HB2	1.77	0.84
1:B:313:LYS:HD3	1:B:317:ASP:OD1	1.79	0.83
1:B:33:PRO:CB	1:B:63:LEU:CD2	2.57	0.83
1:B:420:HIS:ND1	2:B:1450:HEM:CGD	2.41	0.83
1:B:328:ALA:HB3	1:B:329:GLN:HE21	1.40	0.83
1:B:328:ALA:CB	1:B:329:GLN:NE2	2.42	0.83
1:B:115:ALA:O	1:B:123:MET:HG3	1.80	0.82
1:A:308:MET:HE2	1:A:445:PHE:HB3	1.60	0.81
1:B:139:PHE:HD2	1:B:332:TYR:CE2	1.97	0.81
1:B:35:TYR:CE1	1:B:36:PRO:HD2	2.15	0.81
1:B:90:PHE:CD2	1:B:417:ALA:HB3	2.16	0.81
1:B:33:PRO:HB3	1:B:63:LEU:HD23	1.63	0.80
1:B:441:ARG:O	1:B:476:LYS:HD3	1.81	0.80
1:A:113:GLY:HA2	1:A:117:ALA:O	1.80	0.80
1:B:86:HIS:NE2	1:B:413:ILE:HD13	1.97	0.79
1:B:33:PRO:HB2	1:B:63:LEU:CD2	2.13	0.79
1:B:89:PHE:CE2	1:B:90:PHE:CE1	2.69	0.79
1:B:172:GLY:HA2	1:B:297:THR:HG23	1.63	0.79
1:A:339:MET:CE	1:A:437:ALA:HB2	2.13	0.79
1:A:304:MET:O	1:A:308:MET:HG2	1.82	0.78
1:A:247:ARG:NH1	1:A:262:LEU:HD23	1.98	0.78
1:A:359:VAL:C	1:A:360:MET:SD	2.63	0.78
1:A:59:CYS:C	1:A:63:LEU:CD1	2.42	0.77
1:B:51:ASN:ND2	1:B:54:GLU:H	1.83	0.77
1:B:298:ILE:HD13	1:B:463:GLY:CA	2.15	0.77
1:B:313:LYS:CD	1:B:317:ASP:OD1	2.32	0.77
1:B:33:PRO:HB3	1:B:63:LEU:CD2	2.15	0.77
1:A:150:VAL:O	1:A:154:MET:HG3	1.84	0.77
1:B:339:MET:O	1:B:342:ALA:N	2.17	0.76
1:A:240:LEU:HD22	1:A:262:LEU:HD13	1.67	0.76
1:A:331:ASN:H	1:A:334:ASN:ND2	1.82	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:316:LEU:O	1:B:316:LEU:HD12	1.86	0.76
1:B:214:PHE:C	1:B:216:PRO:HD2	2.05	0.76
1:B:175:ILE:HD13	1:B:296:SER:HB2	1.67	0.75
1:B:247:ARG:NH1	1:B:262:LEU:HD23	2.01	0.75
1:B:260:ASP:OD1	1:B:263:GLY:N	2.18	0.75
1:A:211:ALA:CB	1:A:215:MET:HA	2.17	0.74
1:A:76:ARG:HB3	1:A:378:ASP:OD1	1.86	0.74
1:B:191:ARG:HG2	1:B:191:ARG:HH11	1.52	0.74
1:B:348:GLU:OE1	1:B:401:ASP:O	2.04	0.74
1:B:366:GLU:HA	1:B:374:VAL:O	1.88	0.74
1:A:243:ILE:HG22	1:A:247:ARG:HG3	1.70	0.73
1:A:89:PHE:HD2	1:A:90:PHE:CE1	2.07	0.73
1:A:59:CYS:HA	1:A:63:LEU:HD11	1.70	0.73
1:B:314:LYS:HG3	1:B:315:TRP:H	1.54	0.73
1:B:323:ILE:HA	1:B:326:PHE:CD2	2.22	0.73
1:B:211:ALA:O	1:B:215:MET:CB	2.36	0.73
1:B:163:GLY:HA3	1:B:473:TYR:CE2	2.23	0.73
1:A:308:MET:HE2	1:A:445:PHE:HB2	1.67	0.72
1:A:262:LEU:O	1:A:266:LEU:HD13	1.89	0.72
1:B:174:MET:HA	1:B:177:ASN:HD22	1.55	0.72
1:B:319:LEU:HD12	1:B:319:LEU:O	1.89	0.72
1:A:319:LEU:O	1:A:322:GLU:HB3	1.90	0.72
2:A:1450:HEM:CHA	3:A:1460:5PS:H2	2.19	0.72
1:A:351:ARG:O	1:A:351:ARG:HD2	1.90	0.71
1:B:312:ASN:O	1:B:316:LEU:N	2.20	0.71
1:A:297:THR:O	1:A:301:SER:OG	2.08	0.71
1:B:60:LYS:O	1:B:63:LEU:O	2.07	0.71
1:A:306:HIS:O	1:A:312:ASN:ND2	2.18	0.71
1:B:260:ASP:OD1	1:B:262:LEU:N	2.23	0.71
1:B:32:PRO:CA	1:B:372:TYR:CE1	2.74	0.71
1:B:35:TYR:HD1	1:B:36:PRO:HD2	1.44	0.71
1:B:32:PRO:CB	1:B:372:TYR:CD1	2.73	0.71
1:B:98:SER:OG	1:B:101:GLU:OE1	2.05	0.71
1:A:147:GLN:OE1	1:A:330:LEU:CB	2.32	0.71
1:B:326:PHE:CE2	1:B:339:MET:HE2	2.26	0.70
1:A:59:CYS:CA	1:A:63:LEU:CD1	2.70	0.70
1:B:75:GLN:O	1:B:77:VAL:HG23	1.91	0.70
1:B:114:VAL:O	1:B:126:GLN:OE1	2.08	0.70
1:A:277:SER:OG	1:A:280:GLU:HG3	1.92	0.69
1:B:168:LEU:O	1:B:168:LEU:HD23	1.92	0.69
1:B:247:ARG:NH1	1:B:260:ASP:OD2	2.24	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:348:GLU:O	1:B:351:ARG:N	2.25	0.69
1:A:224:PRO:CB	1:A:227:ALA:HB2	2.20	0.69
1:B:442:GLU:O	1:B:476:LYS:CB	2.37	0.69
1:A:211:ALA:HB3	1:A:215:MET:CB	2.23	0.68
1:B:132:GLU:O	1:B:135:THR:HG23	1.93	0.68
1:A:224:PRO:HB2	1:A:227:ALA:CB	2.21	0.68
1:A:59:CYS:CA	1:A:63:LEU:HD11	2.23	0.68
1:B:247:ARG:HH21	1:B:258:THR:N	1.91	0.68
1:A:204:MET:HA	1:A:229:CYS:HB2	1.76	0.68
1:B:66:GLY:O	1:B:81:GLY:N	2.27	0.68
1:B:148:HIS:O	1:B:152:LYS:N	2.25	0.67
1:B:175:ILE:HD12	1:B:296:SER:HB2	1.75	0.67
1:A:401:ASP:OD1	1:A:403:GLU:N	2.25	0.67
1:B:262:LEU:O	1:B:266:LEU:HD22	1.94	0.67
1:A:52:PRO:O	1:A:56:MET:HG3	1.94	0.67
1:B:301:SER:O	1:B:305:LEU:HG	1.94	0.67
1:B:139:PHE:CD2	1:B:332:TYR:HE2	2.06	0.67
2:A:1450:HEM:HBC2	2:A:1450:HEM:HHD	1.77	0.67
1:A:151:ARG:NH2	1:A:328:ALA:O	2.28	0.67
1:B:326:PHE:HE2	1:B:339:MET:HE2	1.60	0.67
1:B:352:ARG:NH1	1:B:398:ARG:NH1	2.42	0.67
1:B:326:PHE:CE2	1:B:339:MET:CE	2.77	0.67
1:B:99:PRO:O	1:B:102:VAL:HG22	1.95	0.67
1:B:361:ARG:NH2	2:B:1450:HEM:CGA	2.52	0.67
1:A:343:GLU:OE1	1:A:433:LYS:NZ	2.28	0.66
1:B:358:MET:HA	1:B:384:PRO:HD2	1.78	0.66
1:B:306:HIS:ND1	1:B:400:TRP:CG	2.62	0.66
2:B:1450:HEM:C1A	3:B:1460:5PS:H2	2.31	0.66
1:B:111:GLY:C	1:B:114:VAL:HG23	2.15	0.66
1:B:325:GLU:OE1	1:B:325:GLU:HA	1.95	0.66
1:A:89:PHE:CD2	1:A:90:PHE:CE1	2.84	0.66
1:B:183:LEU:HD23	1:B:183:LEU:N	2.11	0.66
1:B:33:PRO:HD3	1:B:372:TYR:CE1	2.30	0.66
2:B:1450:HEM:HMC2	2:B:1450:HEM:HBC2	1.78	0.65
1:B:314:LYS:HG3	1:B:315:TRP:N	2.11	0.65
1:A:174:MET:O	1:A:178:THR:OG1	2.12	0.65
1:B:172:GLY:N	1:B:297:THR:HG21	2.00	0.65
1:B:229:CYS:O	1:B:232:ALA:N	2.28	0.65
1:B:59:CYS:O	1:B:63:LEU:HD12	1.96	0.65
1:A:341:PHE:O	1:A:344:ARG:HB2	1.96	0.65
1:A:348:GLU:HA	1:A:348:GLU:OE1	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:351:ARG:O	1:B:388:HIS:ND1	2.29	0.65
1:B:343:GLU:O	1:B:347:ARG:HB2	1.97	0.64
1:A:437:ALA:O	1:A:441:ARG:HB2	1.98	0.64
1:B:154:MET:O	1:B:158:TRP:N	2.30	0.64
1:B:394:PHE:CD2	1:B:404:ARG:HD3	2.33	0.64
1:B:443:TYR:CD1	1:B:475:ARG:HA	2.32	0.64
1:B:211:ALA:O	1:B:213:VAL:O	2.16	0.64
1:B:428:ALA:O	1:B:431:GLN:HB2	1.96	0.64
1:A:358:MET:HA	1:A:384:PRO:HD2	1.79	0.64
1:B:111:GLY:O	1:B:114:VAL:CG2	2.40	0.64
1:B:260:ASP:OD1	1:B:262:LEU:HB3	1.97	0.64
1:B:76:ARG:CB	1:B:378:ASP:OD1	2.45	0.64
1:A:247:ARG:HH12	1:A:262:LEU:HD23	1.62	0.63
1:B:348:GLU:OE2	1:B:351:ARG:NH2	2.32	0.63
1:B:351:ARG:C	1:B:351:ARG:HD2	2.18	0.63
1:A:351:ARG:C	1:A:351:ARG:HD2	2.18	0.63
1:A:211:ALA:O	1:A:215:MET:CA	2.46	0.63
1:B:375:PRO:HD2	1:B:378:ASP:OD2	1.98	0.63
1:A:136:ILE:HA	1:A:139:PHE:CD2	2.34	0.63
1:A:219:LEU:HD12	1:A:219:LEU:N	2.13	0.63
1:B:420:HIS:ND1	2:B:1450:HEM:O2D	2.32	0.63
1:A:305:LEU:HD13	1:A:453:PRO:HD2	1.81	0.63
1:B:135:THR:O	1:B:138:LYS:N	2.32	0.63
1:B:334:ASN:N	1:B:334:ASN:OD1	2.32	0.63
1:A:187:ASP:N	1:A:187:ASP:OD1	2.30	0.62
1:A:188:LEU:HD12	1:A:188:LEU:O	1.98	0.62
1:A:146:ILE:HG13	1:A:434:THR:HG21	1.82	0.62
1:B:215:MET:N	1:B:216:PRO:HD2	2.14	0.62
1:B:201:LEU:HD11	1:B:289:MET:HB3	1.81	0.62
1:B:298:ILE:HD13	1:B:463:GLY:HA3	1.80	0.62
1:B:293:GLN:O	1:B:297:THR:OG1	2.17	0.62
1:A:299:THR:O	1:A:303:SER:OG	2.17	0.61
1:A:272:ASP:OD1	1:A:273:GLY:N	2.33	0.61
1:B:88:ARG:O	1:B:92:PRO:HD3	1.99	0.61
1:A:86:HIS:O	1:A:89:PHE:N	2.32	0.61
2:B:1450:HEM:CHA	3:B:1460:5PS:H2	2.30	0.61
1:B:172:GLY:HA3	1:B:297:THR:HG21	1.75	0.61
1:A:213:VAL:O	1:A:379:ILE:CD1	2.47	0.60
1:B:172:GLY:CA	1:B:297:THR:HG23	2.27	0.60
1:A:229:CYS:O	1:A:232:ALA:N	2.35	0.60
1:A:425:GLN:O	1:A:429:LEU:HD12	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:394:PHE:HD2	1:B:404:ARG:HH11	1.48	0.60
1:A:363:VAL:HG12	1:A:365:ALA:O	2.01	0.60
1:A:158:TRP:HB3	1:A:473:TYR:CZ	2.36	0.60
1:B:247:ARG:CZ	1:B:260:ASP:OD2	2.49	0.60
1:B:44:HIS:CD2	1:B:55:PHE:HZ	2.19	0.60
1:A:229:CYS:SG	1:A:230:ARG:N	2.74	0.60
1:B:344:ARG:HD3	1:B:402:PRO:O	2.01	0.60
1:B:247:ARG:HH12	1:B:262:LEU:HD23	1.67	0.60
1:B:302:TRP:O	1:B:306:HIS:HD2	1.84	0.60
1:B:363:VAL:HG12	1:B:376:LYS:HA	1.82	0.60
1:B:35:TYR:HB3	1:B:44:HIS:HE2	1.67	0.60
1:B:467:ASN:C	1:B:467:ASN:HD22	2.05	0.60
1:A:152:LYS:O	1:A:156:GLU:HG3	2.01	0.59
1:A:51:ASN:C	1:A:51:ASN:HD22	2.06	0.59
2:A:1450:HEM:HBB2	2:A:1450:HEM:HMB2	1.83	0.59
2:B:1450:HEM:C4D	3:B:1460:5PS:H2	2.38	0.59
1:A:243:ILE:HG22	1:A:247:ARG:CG	2.32	0.59
1:B:291:ALA:O	1:B:295:THR:OG1	2.20	0.59
1:B:51:ASN:C	1:B:51:ASN:HD22	2.05	0.59
1:B:66:GLY:C	1:B:81:GLY:HA3	2.11	0.59
1:A:48:PHE:O	1:A:52:PRO:CD	2.48	0.59
1:B:168:LEU:C	1:B:168:LEU:HD23	2.23	0.59
1:B:94:ASN:OD1	1:B:94:ASN:N	2.29	0.59
1:B:328:ALA:CB	1:B:329:GLN:HE22	2.13	0.59
1:A:309:HIS:CE1	1:A:311:LYS:H	2.20	0.59
1:B:110:PHE:O	1:B:114:VAL:HG21	2.03	0.58
1:B:32:PRO:CA	1:B:372:TYR:CD1	2.85	0.58
1:A:146:ILE:HG22	1:A:182:CYS:SG	2.42	0.58
1:A:345:CYS:O	1:A:349:SER:OG	2.21	0.58
1:A:130:LEU:HD23	1:A:423:ILE:HD11	1.85	0.58
1:A:243:ILE:O	1:A:247:ARG:N	2.26	0.58
1:B:191:ARG:HG2	1:B:191:ARG:NH1	2.13	0.58
1:A:331:ASN:HD22	1:A:331:ASN:C	2.07	0.58
1:B:133:GLU:HG3	1:B:261:LEU:HD12	1.86	0.58
1:A:212:ALA:HA	1:A:215:MET:CB	2.33	0.57
1:B:101:GLU:O	1:B:104:THR:HG22	2.04	0.57
1:A:157:ASN:O	1:A:158:TRP:HD1	1.86	0.57
1:A:291:ALA:HB1	3:A:1460:5PS:H3	1.85	0.57
1:B:339:MET:HB3	1:B:342:ALA:HB3	1.85	0.57
1:B:306:HIS:CE1	1:B:400:TRP:CD1	2.92	0.57
1:B:175:ILE:HD13	1:B:296:SER:CB	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:394:PHE:HD2	1:B:404:ARG:HD3	1.69	0.57
1:A:211:ALA:CB	1:A:215:MET:CB	2.82	0.57
1:A:327:PRO:O	1:A:329:GLN:O	2.23	0.57
1:A:59:CYS:SG	1:A:70:ILE:HD11	2.45	0.57
1:A:141:ASN:O	1:A:144:PRO:HD2	2.05	0.56
1:A:385:LEU:O	1:A:389:HIS:ND1	2.33	0.56
1:A:210:PRO:HB3	1:A:460:MET:CE	2.35	0.56
1:B:434:THR:O	1:B:438:THR:OG1	2.23	0.56
1:A:59:CYS:HA	1:A:63:LEU:CD1	2.35	0.56
1:B:197:PHE:O	1:B:200:LEU:HB2	2.06	0.56
1:A:243:ILE:O	1:A:247:ARG:HG3	2.06	0.56
1:B:361:ARG:NH2	2:B:1450:HEM:O2A	2.31	0.56
1:A:89:PHE:HD2	1:A:90:PHE:CD1	2.24	0.56
1:B:244:ILE:O	1:B:248:GLU:N	2.36	0.56
1:B:316:LEU:HD12	1:B:316:LEU:C	2.26	0.56
1:B:429:LEU:HD23	1:B:429:LEU:N	2.21	0.56
1:A:205:GLU:OE2	1:A:293:GLN:NE2	2.38	0.56
1:B:193:ASN:HD22	1:B:193:ASN:C	2.10	0.56
1:A:465:THR:HB	1:A:468:GLN:CG	2.32	0.56
1:B:32:PRO:CB	1:B:372:TYR:CE1	2.89	0.55
2:A:1450:HEM:C4D	3:A:1460:5PS:H2	2.41	0.55
1:B:348:GLU:OE2	1:B:351:ARG:NE	2.38	0.55
1:B:348:GLU:O	1:B:351:ARG:HB3	2.06	0.55
1:A:358:MET:HE2	1:A:383:SER:HB2	1.89	0.55
1:A:243:ILE:HG22	1:A:247:ARG:CD	2.36	0.55
1:B:113:GLY:H	1:B:117:ALA:HB1	1.70	0.55
1:A:214:PHE:O	1:A:214:PHE:HD1	1.90	0.55
1:A:233:ARG:HH11	1:A:237:GLN:HE21	1.52	0.55
1:A:386:LEU:C	1:A:386:LEU:HD12	2.27	0.55
1:A:308:MET:HE3	1:A:445:PHE:HB2	1.85	0.55
1:B:237:GLN:NE2	1:B:279:HIS:ND1	2.55	0.55
1:B:162:GLU:HB3	1:B:474:THR:HG23	1.90	0.54
1:B:175:ILE:CD1	1:B:296:SER:CB	2.83	0.54
1:B:215:MET:N	1:B:216:PRO:CD	2.71	0.54
1:A:204:MET:O	1:A:207:SER:HB3	2.07	0.54
1:A:449:ARG:HG2	1:A:449:ARG:HH11	1.72	0.54
1:B:69:THR:HB	1:B:77:VAL:O	2.07	0.54
1:A:240:LEU:O	1:A:244:ILE:HG13	2.07	0.54
1:A:341:PHE:O	1:A:344:ARG:N	2.41	0.54
1:B:308:MET:SD	1:B:445:PHE:CB	2.96	0.54
1:B:69:THR:HG22	1:B:78:THR:HG23	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:ASN:HD22	1:A:166:ASN:C	2.10	0.54
1:A:211:ALA:O	1:A:215:MET:CB	2.55	0.54
1:A:212:ALA:N	1:A:215:MET:CB	2.70	0.54
1:A:229:CYS:O	1:A:232:ALA:HB3	2.08	0.54
1:A:428:ALA:O	1:A:432:VAL:HG23	2.07	0.54
1:B:170:ASP:OD1	1:B:170:ASP:N	2.41	0.54
1:B:32:PRO:HB3	1:B:372:TYR:CE1	2.41	0.54
1:B:103:TYR:CE2	3:B:1460:5PS:N4	2.74	0.53
1:B:136:ILE:HA	1:B:139:PHE:CE2	2.43	0.53
1:B:152:LYS:O	1:B:156:GLU:HG2	2.08	0.53
1:A:86:HIS:ND1	1:A:410:GLY:O	2.39	0.53
1:A:158:TRP:CZ2	1:A:165:ILE:HD13	2.43	0.53
1:B:315:TRP:HA	1:B:315:TRP:CE3	2.42	0.53
1:B:88:ARG:CB	4:B:2004:HOH:O	2.56	0.53
1:A:158:TRP:HB3	1:A:473:TYR:CE1	2.43	0.53
1:A:233:ARG:NH1	1:A:237:GLN:HE21	2.07	0.53
1:A:277:SER:O	1:A:280:GLU:N	2.40	0.53
1:B:200:LEU:O	1:B:204:MET:HG3	2.09	0.53
1:B:171:CYS:C	1:B:297:THR:CG2	2.77	0.53
1:B:361:ARG:HG2	1:B:361:ARG:NH1	2.24	0.53
1:B:44:HIS:HD2	1:B:55:PHE:HZ	1.56	0.53
1:A:210:PRO:HB3	1:A:460:MET:HE2	1.90	0.53
1:B:308:MET:SD	1:B:445:PHE:HB2	2.49	0.53
1:A:308:MET:HE3	1:A:445:PHE:CB	2.32	0.53
1:B:313:LYS:CA	1:B:316:LEU:HB3	2.33	0.53
1:B:53:LEU:HB2	1:B:385:LEU:HD21	1.90	0.52
1:A:223:LEU:O	1:A:225:GLN:N	2.38	0.52
1:B:104:THR:O	1:B:107:THR:OG1	2.27	0.52
1:A:76:ARG:CB	1:A:378:ASP:OD1	2.57	0.52
1:A:158:TRP:CE2	1:A:165:ILE:HD13	2.45	0.52
1:A:31:LEU:HD21	1:A:375:PRO:HG3	1.92	0.52
1:A:214:PHE:HB2	1:A:217:TRP:HZ3	1.69	0.52
1:B:344:ARG:HH12	1:B:404:ARG:C	2.12	0.52
1:A:85:GLU:OE1	1:A:370:GLY:N	2.43	0.52
1:A:366:GLU:OE2	1:A:373:VAL:HG13	1.98	0.52
1:A:266:LEU:HD12	1:A:266:LEU:N	2.25	0.51
1:B:342:ALA:O	1:B:346:VAL:HG23	2.10	0.51
1:B:412:PHE:CD1	1:B:412:PHE:C	2.83	0.51
1:B:413:ILE:O	1:B:413:ILE:HG13	2.11	0.51
1:B:32:PRO:HA	1:B:372:TYR:CD1	2.46	0.51
1:A:291:ALA:HB1	3:A:1460:5PS:C3	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:319:LEU:HD12	1:B:319:LEU:C	2.30	0.51
1:B:415:PHE:CD1	1:B:425:GLN:HB3	2.46	0.51
1:B:229:CYS:O	1:B:232:ALA:HB3	2.11	0.50
1:B:163:GLY:CA	1:B:473:TYR:CE2	2.92	0.50
1:B:193:ASN:ND2	1:B:196:HIS:HB2	2.26	0.50
1:B:344:ARG:NH1	1:B:404:ARG:O	2.45	0.50
1:B:83:PRO:O	1:B:86:HIS:CB	2.40	0.50
1:B:352:ARG:O	1:B:389:HIS:HE1	1.93	0.50
1:A:386:LEU:HD12	1:A:386:LEU:O	2.11	0.50
1:B:162:GLU:HA	1:B:473:TYR:O	2.10	0.50
1:B:233:ARG:NH1	1:B:282:CYS:SG	2.84	0.50
1:B:35:TYR:CG	1:B:36:PRO:CD	2.93	0.50
1:A:416:GLY:N	2:A:1450:HEM:HMA3	2.27	0.50
1:A:61:ARG:HH11	1:A:61:ARG:CB	2.25	0.50
1:B:362:MET:SD	1:B:363:VAL:N	2.84	0.50
1:A:332:TYR:C	1:A:332:TYR:CD1	2.85	0.50
1:A:32:PRO:HB3	1:A:67:VAL:O	2.11	0.50
1:B:449:ARG:NH1	1:B:451:GLU:HB2	2.27	0.50
1:B:68:PHE:N	1:B:68:PHE:CD1	2.78	0.50
1:A:309:HIS:ND1	1:A:310:PRO:N	2.60	0.50
1:B:224:PRO:O	1:B:227:ALA:HB3	2.11	0.50
1:A:205:GLU:O	1:A:208:LEU:N	2.43	0.50
1:B:110:PHE:C	1:B:114:VAL:HG21	2.32	0.50
1:B:139:PHE:HA	1:B:142:PHE:CG	2.46	0.50
1:B:214:PHE:CA	1:B:216:PRO:HD2	2.41	0.50
1:B:86:HIS:CE1	1:B:413:ILE:HG23	2.47	0.50
1:A:151:ARG:NH1	1:A:442:GLU:OE2	2.39	0.49
1:B:109:VAL:HG12	1:B:286:VAL:HG11	1.94	0.49
1:A:196:HIS:CE1	1:A:235:GLU:OE2	2.65	0.49
1:B:111:GLY:C	1:B:114:VAL:CG2	2.80	0.49
1:A:220:ARG:O	1:B:458:HIS:CD2	2.65	0.49
1:B:59:CYS:O	1:B:63:LEU:CD1	2.60	0.49
1:A:122:ARG:NH2	1:A:270:TYR:CE2	2.80	0.49
1:B:445:PHE:HA	1:B:472:LYS:O	2.13	0.49
1:B:175:ILE:HD12	1:B:297:THR:N	2.27	0.49
1:B:360:MET:C	1:B:361:ARG:HG2	2.32	0.49
1:A:212:ALA:CA	1:A:215:MET:CB	2.91	0.49
1:A:211:ALA:CB	1:A:215:MET:CA	2.89	0.49
1:A:330:LEU:HD22	1:A:334:ASN:HD22	1.78	0.49
1:B:361:ARG:HH11	1:B:361:ARG:HG2	1.77	0.49
1:B:90:PHE:CG	1:B:417:ALA:HB3	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:ILE:HA	1:A:178:THR:HG22	1.95	0.49
1:A:217:TRP:CD1	1:A:217:TRP:C	2.85	0.49
1:A:399:LEU:HD12	1:A:400:TRP:N	2.27	0.49
1:A:166:ASN:HD22	1:A:167:LEU:N	2.11	0.48
1:A:215:MET:N	1:A:216:PRO:CD	2.75	0.48
1:A:85:GLU:OE1	1:A:370:GLY:CA	2.61	0.48
1:B:147:GLN:HA	1:B:150:VAL:HB	1.96	0.48
1:B:110:PHE:C	1:B:233:ARG:HH21	2.02	0.48
1:B:394:PHE:HB3	1:B:404:ARG:NH1	2.29	0.48
1:A:147:GLN:CD	1:A:330:LEU:HB2	2.24	0.48
1:B:201:LEU:CD2	1:B:290:PHE:CD2	2.96	0.48
1:B:33:PRO:HD2	1:B:67:VAL:O	2.14	0.48
1:B:351:ARG:O	1:B:388:HIS:HB3	2.13	0.48
1:B:431:GLN:O	1:B:435:ILE:HG13	2.14	0.48
1:B:174:MET:CA	1:B:177:ASN:HD22	2.21	0.48
1:A:130:LEU:HD21	2:A:1450:HEM:HBC1	1.95	0.48
1:A:423:ILE:HD11	2:A:1450:HEM:HMD2	1.96	0.48
1:A:214:PHE:HB2	1:A:217:TRP:CE3	2.45	0.48
1:B:315:TRP:HE3	1:B:315:TRP:HA	1.78	0.48
1:B:319:LEU:HD12	1:B:323:ILE:HG23	1.96	0.48
1:B:364:LYS:O	1:B:376:LYS:HD2	2.14	0.48
1:B:359:VAL:C	1:B:360:MET:HE2	2.34	0.48
1:A:219:LEU:CD1	1:A:219:LEU:N	2.77	0.48
1:B:188:LEU:HA	1:B:243:ILE:CD1	2.44	0.48
1:B:260:ASP:OD1	1:B:262:LEU:CA	2.61	0.48
1:A:122:ARG:NH2	1:A:270:TYR:OH	2.46	0.48
1:B:86:HIS:CE1	1:B:411:ALA:O	2.67	0.48
1:B:70:ILE:O	1:B:77:VAL:HB	2.13	0.47
1:B:90:PHE:N	1:B:90:PHE:CD1	2.82	0.47
1:B:331:ASN:HD22	1:B:331:ASN:C	2.18	0.47
1:A:214:PHE:CD1	1:A:214:PHE:C	2.86	0.47
1:A:449:ARG:HG2	1:A:449:ARG:NH1	2.30	0.47
2:B:1450:HEM:CMC	2:B:1450:HEM:HBC2	2.41	0.47
1:B:109:VAL:HG13	1:B:204:MET:SD	2.54	0.47
1:A:170:ASP:N	1:A:170:ASP:OD1	2.46	0.47
1:B:308:MET:SD	1:B:445:PHE:HB3	2.54	0.47
1:B:260:ASP:OD1	1:B:262:LEU:CB	2.63	0.47
1:B:83:PRO:C	1:B:86:HIS:HB2	2.27	0.47
1:A:166:ASN:OD1	1:A:466:LEU:HD11	2.15	0.47
1:A:211:ALA:O	1:A:214:PHE:N	2.48	0.47
1:B:166:ASN:C	1:B:166:ASN:HD22	2.18	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:152:LYS:O	1:B:156:GLU:CG	2.63	0.47
1:A:238:LYS:O	1:A:242:GLU:HG3	2.14	0.47
1:A:344:ARG:NH1	1:A:405:ASP:OD1	2.45	0.47
1:B:306:HIS:ND1	1:B:400:TRP:CD1	2.82	0.47
1:B:139:PHE:CD2	1:B:332:TYR:CE2	2.89	0.47
1:B:359:VAL:CA	1:B:360:MET:HE2	2.45	0.47
3:A:1460:5PS:O2	3:A:1460:5PS:H4	2.14	0.46
1:B:345:CYS:O	1:B:349:SER:OG	2.24	0.46
1:A:215:MET:N	1:A:216:PRO:HD3	2.30	0.46
1:A:214:PHE:C	1:A:216:PRO:CD	2.84	0.46
1:A:277:SER:O	1:A:279:HIS:N	2.48	0.46
1:A:44:HIS:HD2	1:A:70:ILE:HG13	1.80	0.46
1:A:260:ASP:OD1	1:A:260:ASP:N	2.48	0.46
1:A:315:TRP:CE2	1:A:402:PRO:HD2	2.51	0.46
1:B:171:CYS:C	1:B:297:THR:HG23	2.36	0.46
1:B:412:PHE:CE1	1:B:414:GLY:N	2.81	0.46
1:A:425:GLN:HG3	1:A:426:LYS:N	2.30	0.46
1:A:157:ASN:C	1:A:158:TRP:CD1	2.89	0.46
1:A:61:ARG:HB2	1:A:61:ARG:HH11	1.81	0.46
1:B:313:LYS:HD2	1:B:317:ASP:OD1	2.13	0.46
1:B:44:HIS:HD2	1:B:55:PHE:CZ	2.34	0.46
1:A:200:LEU:HA	1:A:200:LEU:HD23	1.77	0.46
1:A:44:HIS:HB2	1:A:71:SER:O	2.16	0.46
1:B:394:PHE:HD2	1:B:404:ARG:NH1	2.13	0.46
1:B:396:ASN:ND2	1:B:399:LEU:HB2	2.31	0.46
1:B:466:LEU:O	1:B:466:LEU:HD12	2.16	0.46
1:A:449:ARG:CD	1:A:468:GLN:OE1	2.60	0.45
1:A:90:PHE:O	1:A:91:SER:CB	2.64	0.45
1:B:44:HIS:N	1:B:44:HIS:ND1	2.64	0.45
1:A:68:PHE:N	1:A:68:PHE:CD1	2.83	0.45
1:A:92:PRO:HG2	1:A:97:LEU:HD12	1.98	0.45
1:A:99:PRO:HB2	1:A:103:TYR:CD2	2.51	0.45
1:A:409:ASP:OD1	1:A:409:ASP:N	2.49	0.45
1:B:183:LEU:N	1:B:183:LEU:CD2	2.78	0.45
1:A:154:MET:O	1:A:158:TRP:N	2.45	0.45
1:B:96:ILE:O	1:B:364:LYS:HB2	2.16	0.45
1:A:162:GLU:HA	1:A:473:TYR:O	2.16	0.45
1:B:86:HIS:CE1	1:B:413:ILE:CD1	2.94	0.45
1:A:233:ARG:NH1	1:A:237:GLN:NE2	2.65	0.45
1:B:120:TYR:N	1:B:121:PRO:CD	2.80	0.45
1:B:52:PRO:O	1:B:55:PHE:HB3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:MET:HG3	1:A:381:ALA:HB2	1.98	0.45
1:A:315:TRP:CE3	1:A:315:TRP:HA	2.51	0.45
1:B:344:ARG:CD	1:B:402:PRO:O	2.65	0.45
1:B:327:PRO:C	1:B:441:ARG:HH12	2.18	0.45
1:A:106:MET:HE2	1:A:110:PHE:CZ	2.52	0.45
1:A:201:LEU:HD11	1:A:289:MET:HB3	1.99	0.45
1:A:233:ARG:NH1	1:A:282:CYS:SG	2.90	0.45
1:A:83:PRO:HA	1:A:86:HIS:CE1	2.51	0.45
1:B:360:MET:O	1:B:361:ARG:HG2	2.17	0.45
1:A:153:PHE:CD1	1:A:153:PHE:C	2.91	0.44
1:B:389:HIS:HA	1:B:397:PRO:HB3	1.98	0.44
1:B:326:PHE:C	1:B:441:ARG:HH22	2.13	0.44
1:A:470:LEU:O	1:A:471:VAL:HG23	2.18	0.44
1:B:193:ASN:HD21	1:B:196:HIS:HB2	1.81	0.44
1:B:352:ARG:HH11	1:B:398:ARG:NH1	2.14	0.44
1:A:32:PRO:HA	1:A:372:TYR:CD1	2.52	0.44
1:B:113:GLY:N	1:B:117:ALA:CB	2.80	0.44
1:B:302:TRP:HA	1:B:305:LEU:HD12	1.98	0.44
1:B:51:ASN:HD22	1:B:53:LEU:N	2.15	0.44
1:A:115:ALA:O	1:A:123:MET:HG3	2.17	0.44
1:A:214:PHE:CB	1:A:217:TRP:CZ3	2.88	0.44
1:A:72:ILE:HD13	1:A:72:ILE:HA	1.82	0.44
1:B:319:LEU:HD11	1:B:323:ILE:CG2	2.47	0.44
1:B:56:MET:HB2	1:B:386:LEU:HD13	2.00	0.44
1:B:90:PHE:HD2	1:B:417:ALA:O	2.01	0.44
1:A:196:HIS:HE1	1:A:235:GLU:OE2	2.01	0.44
1:A:143:VAL:CG2	1:A:332:TYR:HA	2.39	0.44
1:B:237:GLN:HG3	1:B:282:CYS:SG	2.58	0.44
1:A:110:PHE:C	1:A:233:ARG:HH21	1.96	0.44
1:B:107:THR:HG22	1:B:117:ALA:HB2	2.00	0.44
1:B:107:THR:CG2	1:B:117:ALA:HB2	2.48	0.44
1:B:148:HIS:CD2	1:B:148:HIS:C	2.90	0.44
1:B:424:GLY:HA3	2:B:1450:HEM:C4C	2.53	0.44
2:B:1450:HEM:HBB2	2:B:1450:HEM:HMB2	2.00	0.44
1:B:323:ILE:CA	1:B:326:PHE:CD2	2.99	0.44
1:B:86:HIS:NE2	1:B:413:ILE:CD1	2.77	0.44
1:A:60:LYS:HG3	1:A:66:GLY:HA3	1.99	0.43
1:B:192:LEU:HA	1:B:192:LEU:HD12	1.74	0.43
1:A:469:CYS:O	1:A:471:VAL:HG23	2.18	0.43
1:B:102:VAL:HG23	1:B:103:TYR:N	2.33	0.43
1:A:248:GLU:OE1	1:A:248:GLU:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:291:ALA:O	1:A:295:THR:OG1	2.36	0.43
1:A:398:ARG:HD2	1:A:398:ARG:HA	1.75	0.43
1:A:307:LEU:HD23	1:A:307:LEU:HA	1.87	0.43
1:A:352:ARG:HH12	1:A:398:ARG:CZ	2.31	0.43
1:B:146:ILE:C	1:B:150:VAL:HG23	2.38	0.43
1:B:186:GLU:O	1:B:190:LYS:N	2.32	0.43
1:B:277:SER:HG	1:B:280:GLU:HG3	1.82	0.43
1:B:330:LEU:HD22	1:B:334:ASN:HB2	2.01	0.43
1:B:69:THR:HA	1:B:78:THR:HA	2.01	0.43
1:B:193:ASN:O	1:B:196:HIS:HB3	2.18	0.43
2:A:1450:HEM:C1A	3:A:1460:5PS:C2	2.91	0.43
1:A:121:PRO:O	1:A:125:GLU:HG3	2.19	0.43
1:A:130:LEU:HD21	2:A:1450:HEM:CBC	2.49	0.43
1:A:157:ASN:ND2	1:A:157:ASN:N	2.67	0.43
1:B:166:ASN:HD22	1:B:169:GLU:H	1.66	0.43
1:B:360:MET:N	1:B:360:MET:CE	2.82	0.43
1:A:277:SER:HG	1:A:280:GLU:HG3	1.84	0.42
1:B:328:ALA:HB1	1:B:329:GLN:HE22	1.81	0.42
1:B:358:MET:HE2	1:B:358:MET:HB3	1.69	0.42
1:A:393:ALA:HB2	1:A:408:VAL:HG23	2.01	0.42
1:B:162:GLU:CB	1:B:474:THR:HG23	2.49	0.42
1:B:166:ASN:ND2	1:B:169:GLU:H	2.17	0.42
1:B:323:ILE:O	1:B:326:PHE:CD2	2.73	0.42
1:B:330:LEU:HD23	1:B:330:LEU:HA	1.91	0.42
1:B:344:ARG:NH1	1:B:404:ARG:C	2.72	0.42
1:A:358:MET:HB3	1:A:358:MET:HE3	1.88	0.42
1:B:139:PHE:N	1:B:139:PHE:CD1	2.83	0.42
1:B:329:GLN:N	1:B:329:GLN:NE2	2.67	0.42
1:A:351:ARG:C	1:A:351:ARG:CD	2.85	0.42
1:A:166:ASN:OD1	1:A:466:LEU:CD1	2.67	0.42
1:B:292:GLY:O	1:B:296:SER:OG	2.22	0.42
1:B:69:THR:HG22	1:B:78:THR:OG1	2.19	0.42
1:A:295:THR:HB	2:A:1450:HEM:CAB	2.49	0.42
1:A:214:PHE:CD1	1:A:214:PHE:O	2.70	0.42
1:B:127:LEU:O	1:B:130:LEU:N	2.52	0.42
1:B:233:ARG:HH12	1:B:282:CYS:HB3	1.84	0.42
1:B:201:LEU:HD22	1:B:290:PHE:CD2	2.55	0.42
1:B:421:LYS:O	1:B:422:CYS:C	2.54	0.42
1:B:166:ASN:C	1:B:166:ASN:ND2	2.72	0.42
1:B:383:SER:HA	1:B:384:PRO:HD2	1.92	0.42
1:A:293:GLN:HG3	1:A:294:HIS:N	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:ILE:HD12	1:A:461:VAL:HG12	2.02	0.42
1:A:359:VAL:O	1:A:381:ALA:HA	2.19	0.42
1:B:295:THR:HB	2:B:1450:HEM:CAB	2.50	0.42
1:B:363:VAL:CG1	1:B:376:LYS:HA	2.47	0.42
1:A:100:ARG:HA	1:A:116:TYR:HB3	2.01	0.42
3:A:1460:5PS:C4	3:A:1460:5PS:O2	2.67	0.42
1:A:439:ALA:O	1:A:443:TYR:HB2	2.20	0.42
1:B:306:HIS:HE1	1:B:352:ARG:HE	1.66	0.42
1:B:329:GLN:N	1:B:329:GLN:CD	2.73	0.42
1:B:348:GLU:OE1	1:B:401:ASP:C	2.58	0.42
1:B:89:PHE:HE2	1:B:90:PHE:CZ	2.37	0.42
1:A:61:ARG:HH11	1:A:61:ARG:CG	2.32	0.41
1:B:394:PHE:CD2	1:B:404:ARG:NH1	2.85	0.41
1:B:443:TYR:CE1	1:B:475:ARG:HB2	2.56	0.41
1:B:51:ASN:HD21	1:B:53:LEU:HB3	1.85	0.41
1:B:118:ALA:HB1	1:B:122:ARG:HG2	2.03	0.41
1:B:33:PRO:HD3	1:B:372:TYR:HE1	1.84	0.41
1:B:384:PRO:O	1:B:388:HIS:CD2	2.73	0.41
1:A:188:LEU:HD12	1:A:188:LEU:C	2.41	0.41
1:A:334:ASN:HA	1:A:338:GLU:HB2	2.00	0.41
1:B:326:PHE:C	1:B:441:ARG:NH2	2.62	0.41
1:A:101:GLU:O	1:A:104:THR:HG22	2.20	0.41
1:A:120:TYR:N	1:A:121:PRO:HD2	2.36	0.41
1:A:339:MET:HE3	1:A:437:ALA:HB2	1.99	0.41
1:B:113:GLY:CA	1:B:117:ALA:HB3	2.51	0.41
1:A:278:LEU:HA	1:A:281:VAL:HB	2.03	0.41
1:A:331:ASN:ND2	1:A:334:ASN:ND2	2.69	0.41
1:A:418:GLY:O	1:A:419:VAL:C	2.57	0.41
1:B:86:HIS:CE1	1:B:413:ILE:CG2	3.04	0.41
1:B:211:ALA:O	1:B:215:MET:N	2.53	0.41
1:B:298:ILE:HD12	1:B:461:VAL:HG12	2.03	0.41
1:A:75:GLN:HG2	1:A:217:TRP:CZ3	2.55	0.41
1:B:315:TRP:NE1	1:B:402:PRO:HD2	2.35	0.41
1:B:394:PHE:CB	1:B:404:ARG:NH1	2.84	0.41
1:B:51:ASN:ND2	1:B:53:LEU:N	2.69	0.41
1:A:136:ILE:HA	1:A:139:PHE:CE2	2.56	0.41
1:A:330:LEU:HA	1:A:330:LEU:HD23	1.79	0.41
2:B:1450:HEM:C1A	3:B:1460:5PS:C2	3.02	0.41
1:B:298:ILE:HD13	1:B:463:GLY:HA2	2.02	0.41
1:B:398:ARG:HA	1:B:398:ARG:HD2	1.89	0.41
1:B:147:GLN:O	1:B:151:ARG:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1450:HEM:C4D	3:A:1460:5PS:C2	3.04	0.41
1:A:453:PRO:HB3	1:A:468:GLN:HB2	2.03	0.41
1:B:89:PHE:CE2	1:B:90:PHE:CZ	3.09	0.41
1:A:320:HIS:O	1:A:324:ASP:N	2.55	0.40
1:B:183:LEU:O	1:B:184:PHE:CD1	2.74	0.40
1:A:270:TYR:OH	1:A:280:GLU:OE1	2.35	0.40
1:A:295:THR:HG22	1:A:355:PRO:HB2	2.03	0.40
1:A:331:ASN:H	1:A:334:ASN:HD22	1.65	0.40
1:A:423:ILE:CD1	2:A:1450:HEM:HMD2	2.51	0.40
1:B:102:VAL:HG23	1:B:103:TYR:HD1	1.74	0.40
1:B:467:ASN:C	1:B:467:ASN:ND2	2.73	0.40
1:A:168:LEU:HA	1:A:168:LEU:HD12	1.84	0.40
1:A:277:SER:O	1:A:278:LEU:C	2.59	0.40
1:A:348:GLU:HG2	1:A:401:ASP:O	2.22	0.40
1:B:473:TYR:CD1	1:B:473:TYR:C	2.94	0.40
1:B:361:ARG:O	1:B:380:ILE:HG22	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	437/467 (94%)	414 (95%)	21 (5%)	2 (0%)	32	71
1	B	438/467 (94%)	427 (98%)	9 (2%)	2 (0%)	32	71
All	All	875/934 (94%)	841 (96%)	30 (3%)	4 (0%)	32	71

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	91	SER
1	A	278	LEU

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Mol	Chain	Res	Type
1	B	222	PRO
1	B	339	MET

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	343/408 (84%)	289 (84%)	54 (16%)	3	13
1	B	314/408 (77%)	253 (81%)	61 (19%)	1	7
All	All	657/816 (80%)	542 (82%)	115 (18%)	2	10

All (115) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	THR
1	A	39	VAL
1	A	51	ASN
1	A	61	ARG
1	A	63	LEU
1	A	65	SER
1	A	68	PHE
1	A	87	SER
1	A	88	ARG
1	A	90	PHE
1	A	104	THR
1	A	138	LYS
1	A	146	ILE
1	A	162	GLU
1	A	166	ASN
1	A	170	ASP
1	A	178	THR
1	A	188	LEU
1	A	192	LEU
1	A	197	PHE
1	A	203	LYS

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Mol	Chain	Res	Type
1	A	205	GLU
1	A	213	VAL
1	A	214	PHE
1	A	217	TRP
1	A	230	ARG
1	A	251	GLU
1	A	274	THR
1	A	278	LEU
1	A	282	CYS
1	A	301	SER
1	A	303	SER
1	A	321	LYS
1	A	331	ASN
1	A	332	TYR
1	A	336	MET
1	A	349	SER
1	A	352	ARG
1	A	353	ASP
1	A	360	MET
1	A	366	GLU
1	A	378	ASP
1	A	386	LEU
1	A	390	ASP
1	A	391	GLU
1	A	403	GLU
1	A	406	GLU
1	A	409	ASP
1	A	442	GLU
1	A	444	ASP
1	A	446	GLN
1	A	459	THR
1	A	467	ASN
1	A	470	LEU
1	B	34	VAL
1	B	39	VAL
1	B	44	HIS
1	B	51	ASN
1	B	53	LEU
1	B	63	LEU
1	B	86	HIS
1	B	94	ASN
1	B	106	MET

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Mol	Chain	Res	Type
1	B	107	THR
1	B	122	ARG
1	B	139	PHE
1	B	154	MET
1	B	156	GLU
1	B	166	ASN
1	B	170	ASP
1	B	183	LEU
1	B	191	ARG
1	B	193	ASN
1	B	199	GLN
1	B	203	LYS
1	B	205	GLU
1	B	229	CYS
1	B	231	GLU
1	B	233	ARG
1	B	242	GLU
1	B	266	LEU
1	B	274	THR
1	B	290	PHE
1	B	295	THR
1	B	296	SER
1	B	297	THR
1	B	303	SER
1	B	304	MET
1	B	307	LEU
1	B	314	LYS
1	B	316	LEU
1	B	319	LEU
1	B	320	HIS
1	B	321	LYS
1	B	323	ILE
1	B	325	GLU
1	B	331	ASN
1	B	333	ASP
1	B	334	ASN
1	B	336	MET
1	B	339	MET
1	B	348	GLU
1	B	356	LEU
1	B	360	MET
1	B	361	ARG

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Mol	Chain	Res	Type
1	B	366	GLU
1	B	402	PRO
1	B	438	THR
1	B	449	ARG
1	B	458	HIS
1	B	459	THR
1	B	460	MET
1	B	466	LEU
1	B	467	ASN
1	B	474	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	166	ASN
1	A	196	HIS
1	A	237	GLN
1	A	294	HIS
1	A	306	HIS
1	A	331	ASN
1	A	334	ASN
1	B	51	ASN
1	B	148	HIS
1	B	166	ASN
1	B	193	ASN
1	B	237	GLN
1	B	293	GLN
1	B	329	GLN
1	B	331	ASN
1	B	431	GLN
1	B	458	HIS
1	B	467	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

4 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$
2	HEM	A	1450	1,3	28,50,50	1.24	5 (17%)	17,82,82	1.75	3 (17%)
3	5PS	A	1460	2	40,41,41	1.36	3 (7%)	51,57,57	1.39	6 (11%)
2	HEM	B	1450	1,3	28,50,50	1.21	3 (10%)	17,82,82	1.63	2 (11%)
3	5PS	B	1460	2	40,41,41	1.38	3 (7%)	51,57,57	1.55	8 (15%)

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	HEM	A	1450	1,3	-	0/6/54/54	0/0/8/8
3	5PS	A	1460	2	-	0/23/24/24	0/5/5/5
2	HEM	B	1450	1,3	-	0/6/54/54	0/0/8/8
3	5PS	B	1460	2	-	0/23/24/24	0/5/5/5

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1450	HEM	C1B-NB	-3.64	1.32	1.36
3	B	1460	5PS	C5-N2	-3.45	1.34	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1460	5PS	C5-N2	-3.15	1.35	1.41
3	A	1460	5PS	C25-C24	-2.74	1.37	1.41
2	B	1450	HEM	C4C-NC	-2.42	1.33	1.36
2	A	1450	HEM	C4B-NB	-2.37	1.31	1.36
2	A	1450	HEM	C3C-C2C	-2.35	1.37	1.40
3	B	1460	5PS	C28-C29	-2.29	1.37	1.42
2	A	1450	HEM	C4C-NC	-2.25	1.34	1.36
2	B	1450	HEM	C1C-NC	-2.16	1.34	1.36
2	A	1450	HEM	C1D-ND	-2.15	1.31	1.36
2	A	1450	HEM	C1B-NB	-2.01	1.34	1.36
3	A	1460	5PS	C9-C20	5.77	1.47	1.38
3	B	1460	5PS	C9-C20	5.79	1.47	1.38

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1460	5PS	C19-C20-C9	-3.75	119.82	123.62
3	A	1460	5PS	C19-C20-C9	-3.64	119.93	123.62
3	B	1460	5PS	C5-N2-C6	-3.40	120.65	127.29
2	A	1450	HEM	CAD-CBD-CGD	-3.30	107.02	112.66
2	A	1450	HEM	CBA-CAA-C2A	-3.29	106.19	112.48
3	B	1460	5PS	O1-C8-C9	-3.01	115.30	120.98
3	A	1460	5PS	C5-N2-C6	-2.83	121.76	127.29
3	A	1460	5PS	O2-C6-C7	-2.67	114.67	120.43
3	B	1460	5PS	C27-C28-C29	-2.53	117.26	120.88
2	B	1450	HEM	CAD-C3D-C2D	-2.15	122.87	129.00
3	A	1460	5PS	C27-C28-C29	-2.10	117.88	120.88
3	B	1460	5PS	C11-C12-C19	2.42	121.29	118.16
3	A	1460	5PS	C9-C8-N3	2.60	121.82	116.88
3	B	1460	5PS	C10-C9-C20	2.66	119.74	116.64
3	B	1460	5PS	C9-C8-N3	2.91	122.41	116.88
3	A	1460	5PS	C28-C29-C24	3.20	122.41	118.17
3	B	1460	5PS	C28-C29-C24	3.48	122.79	118.17
2	A	1450	HEM	CBD-CAD-C3D	4.11	120.31	112.47
2	B	1450	HEM	CBD-CAD-C3D	4.23	120.54	112.47

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 34 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1450	HEM	13	0
3	A	1460	5PS	9	0
2	B	1450	HEM	16	0
3	B	1460	5PS	5	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	441/467 (94%)	-0.48	1 (0%) 94 89	43, 70, 99, 121	0
1	B	442/467 (94%)	-0.34	1 (0%) 94 89	59, 91, 120, 135	0
All	All	883/934 (94%)	-0.41	2 (0%) 94 89	43, 79, 116, 135	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	373	VAL	2.0
1	B	436	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	5PS	B	1460	37/37	0.93	0.34	3.16	69,74,95,109	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	HEM	B	1450	43/43	0.97	0.25	1.31	61,70,80,80	0
3	5PS	A	1460	37/37	0.94	0.26	1.14	54,63,78,100	0
2	HEM	A	1450	43/43	0.98	0.20	-0.02	40,48,55,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.