



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:54 PM UTC

PDB ID : 4COX / pdb_00004cox
Title : CYCLOOXYGENASE-2 (PROSTAGLANDIN SYNTHASE-2) COM-
PLEXED WITH A NON-SELECTIVE INHIBITOR, INDOMETHACIN
Authors : Kurumbail, R.; Stallings, W.
Deposited on : 1996-12-18
Resolution : 2.90 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

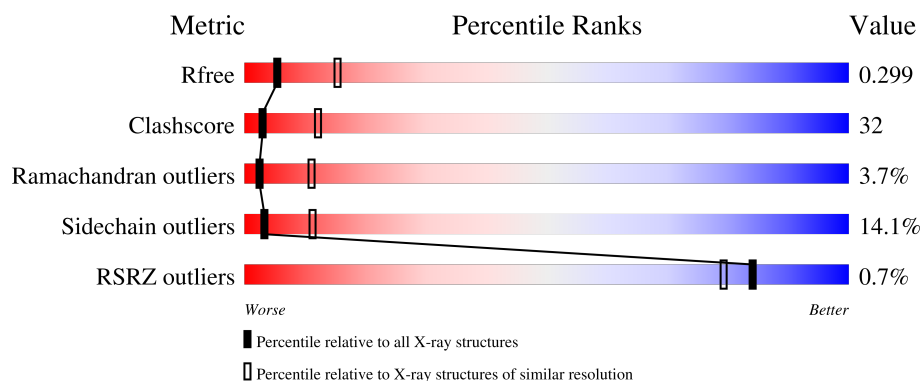
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	587	% 38% 41% 14% • 6%
1	B	587	39% 42% 12% • 6%
1	C	587	% 38% 42% 12% • 6%
1	D	587	% 38% 43% 13% • 6%

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
4	IMN	B	701	-	-	X	-
4	IMN	D	701	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 18332 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.

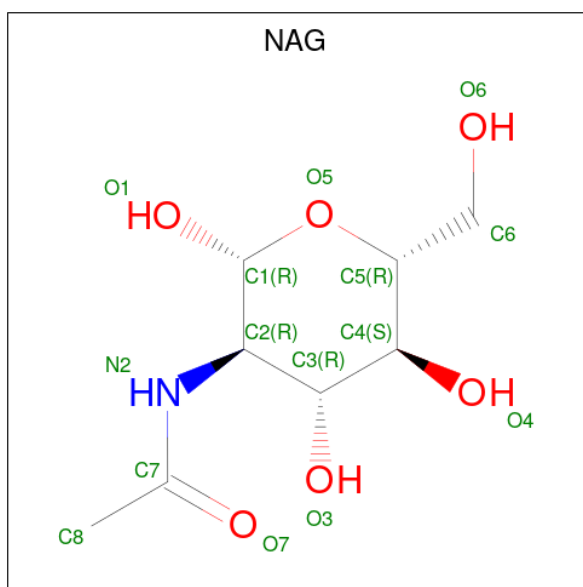
- Molecule 1 is a protein called CYCLOOXYGENASE-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	552	Total	C	N	O	S	0	0	0
			4473	2886	748	814	25			
1	B	552	Total	C	N	O	S	0	0	0
			4473	2886	748	814	25			
1	C	552	Total	C	N	O	S	0	0	0
			4473	2886	748	814	25			
1	D	552	Total	C	N	O	S	0	0	0
			4473	2886	748	814	25			

There are 8 discrepancies between the modelled and reference sequences:

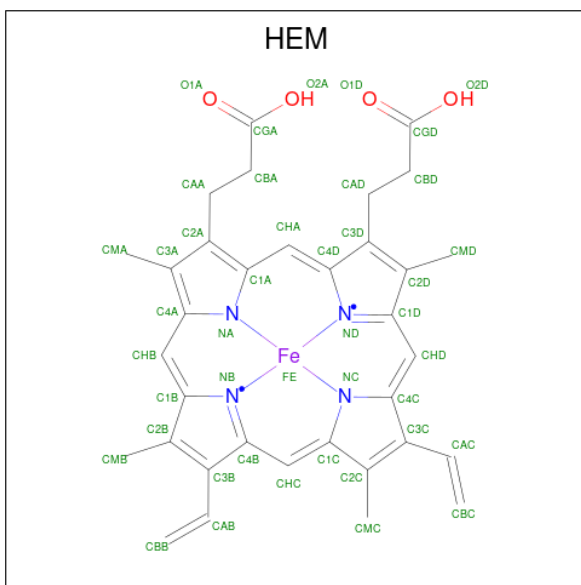
Chain	Residue	Modelled	Actual	Comment	Reference
A	310	GLN	ASN	conflict	UNP Q05769
A	333	LYS	ARG	conflict	UNP Q05769
B	310	GLN	ASN	conflict	UNP Q05769
B	333	LYS	ARG	conflict	UNP Q05769
C	310	GLN	ASN	conflict	UNP Q05769
C	333	LYS	ARG	conflict	UNP Q05769
D	310	GLN	ASN	conflict	UNP Q05769
D	333	LYS	ARG	conflict	UNP Q05769

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



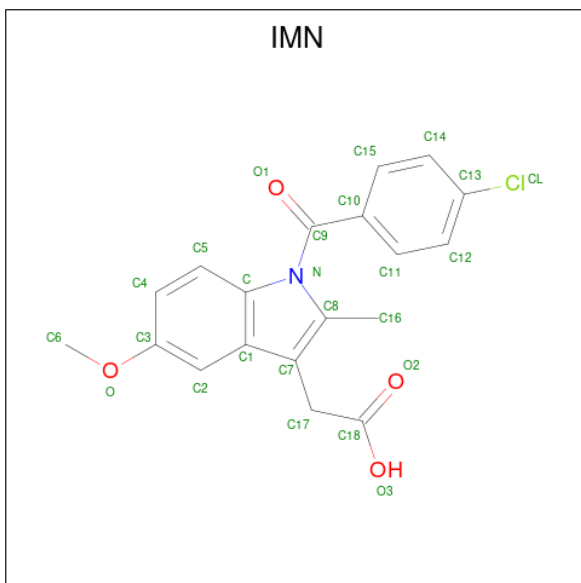
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	C	1	Total	C	N	O	0	0
			14	8	1	5		
2	C	1	Total	C	N	O	0	0
			14	8	1	5		
2	C	1	Total	C	N	O	0	0
			14	8	1	5		
2	D	1	Total	C	N	O	0	0
			14	8	1	5		
2	D	1	Total	C	N	O	0	0
			14	8	1	5		
2	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 4 is INDOMETHACIN (CCD ID: IMN) (formula: $C_{19}H_{16}ClNO_4$).

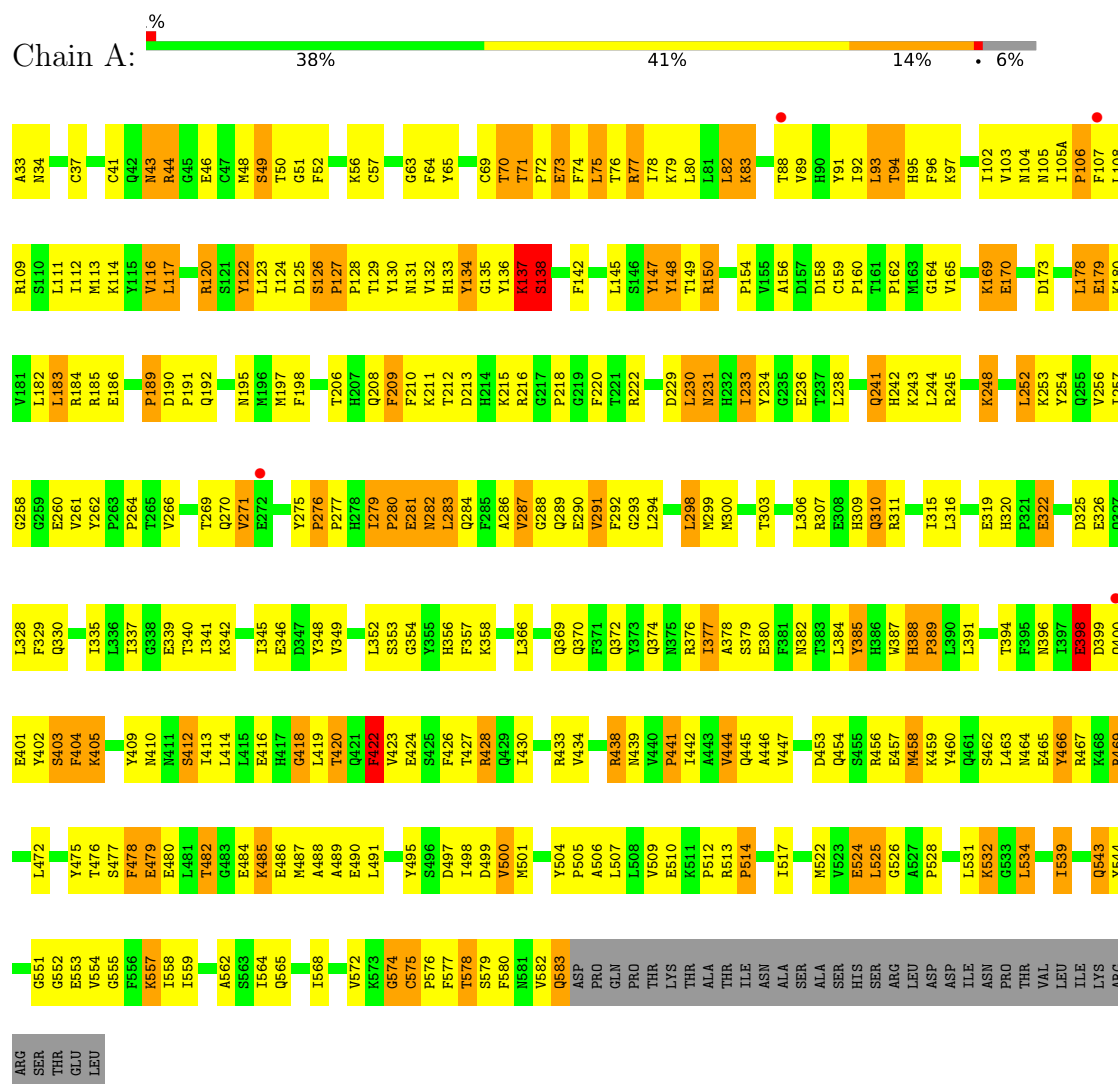


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 25	C 19	Cl 1	N 1	O 4	0	0
4	B	1	Total 25	C 19	Cl 1	N 1	O 4	0	0
4	C	1	Total 25	C 19	Cl 1	N 1	O 4	0	0
4	D	1	Total 25	C 19	Cl 1	N 1	O 4	0	0

3 Residue-property plots

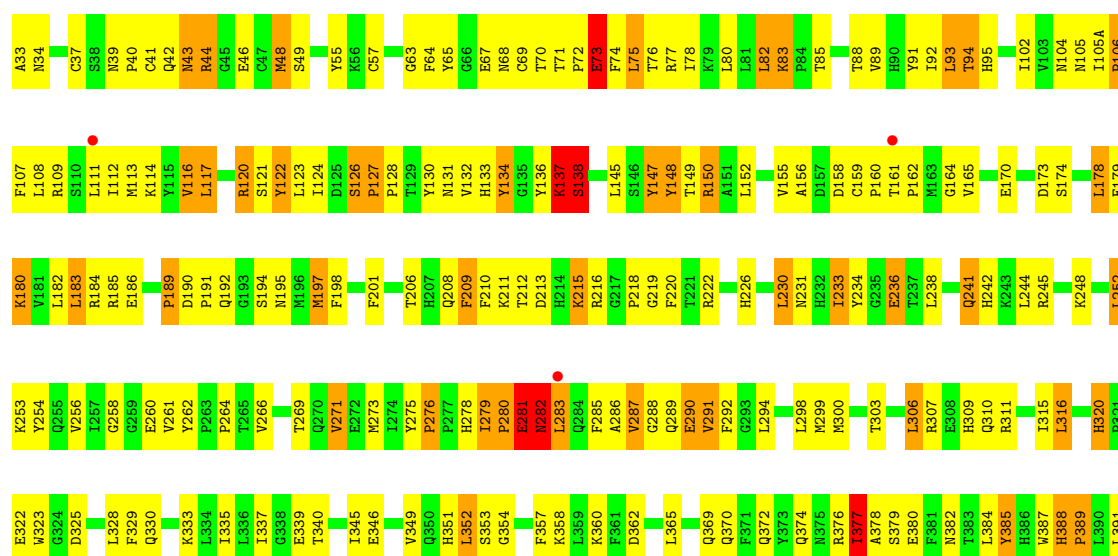
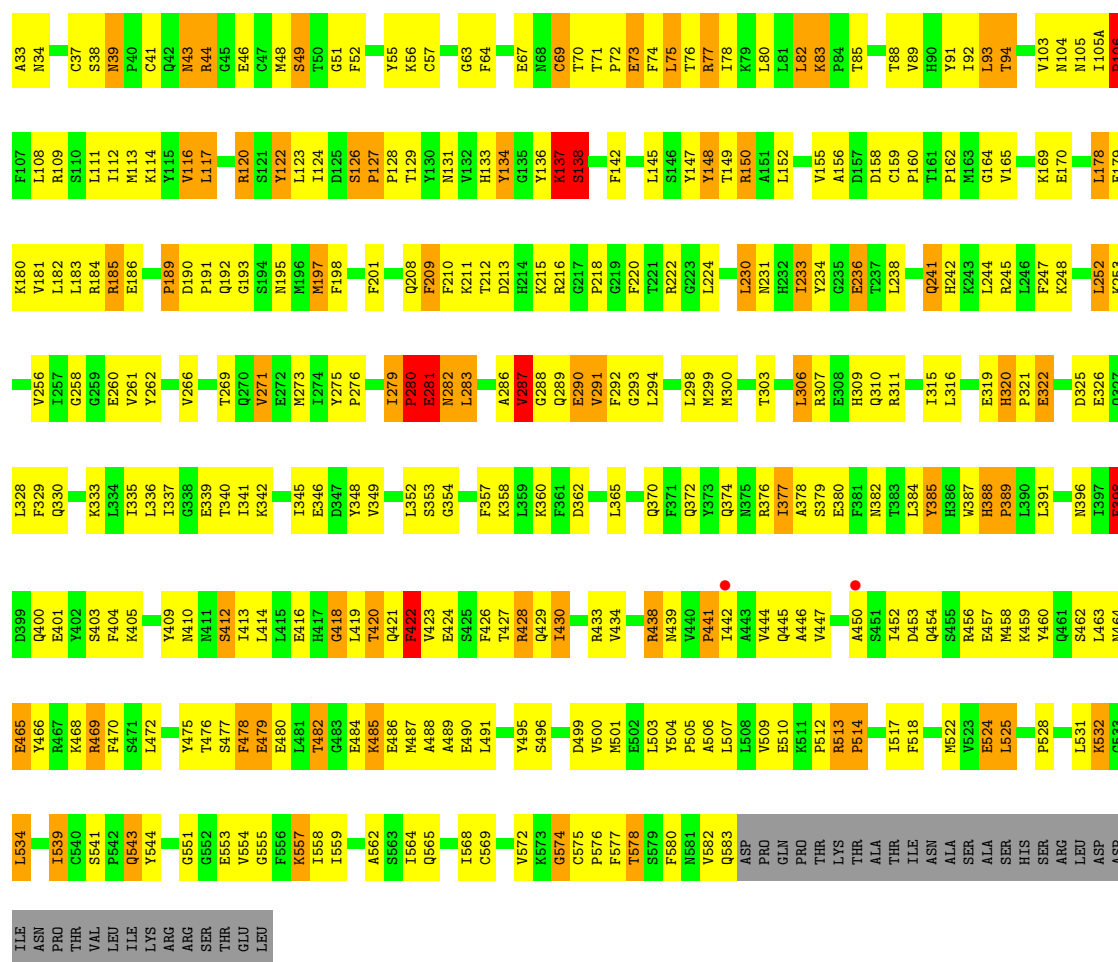
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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.

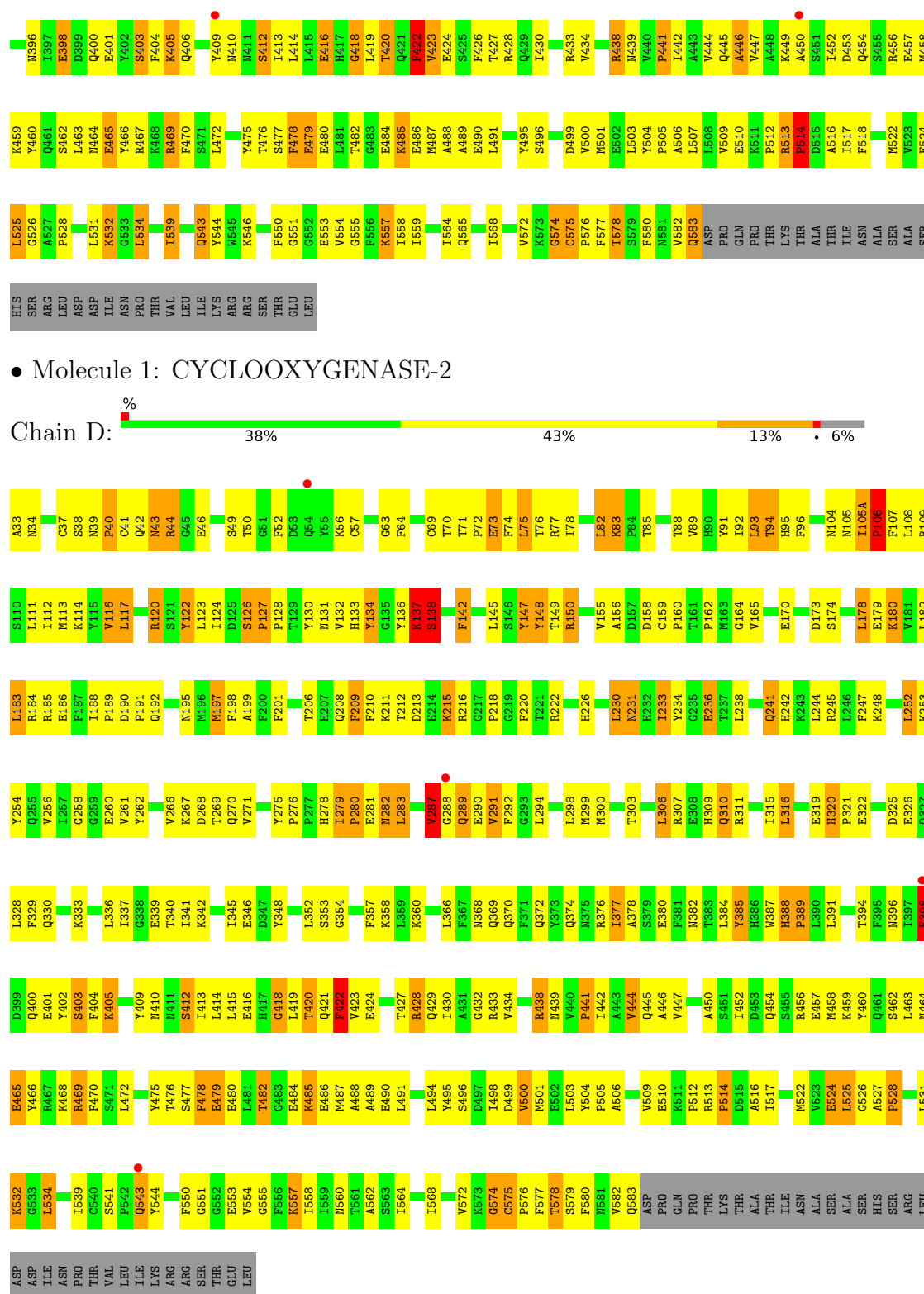
• Molecule 1: CYCLOOXYGENASE-2



• Molecule 1: CYCLOOXYGENASE-2







4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	179.80Å 133.60Å 118.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.90 8.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	68.0 (8.00-2.90) 72.2 (8.00-2.90)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.88Å)	Xtrriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.219 , 0.309 0.228 , 0.299	Depositor DCC
R_{free} test set	4858 reflections (10.08%)	wwPDB-VP
Wilson B-factor (Å ²)	31.8	Xtrriage
Anisotropy	0.735	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 70.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	18332	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 30.54 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2860e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

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

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	0.84	2/4600 (0.0%)	1.23	45/6237 (0.7%)
1	B	0.84	3/4600 (0.1%)	1.24	46/6237 (0.7%)
1	C	0.81	0/4600	1.24	50/6237 (0.8%)
1	D	0.81	0/4600	1.24	52/6237 (0.8%)
All	All	0.82	5/18400 (0.0%)	1.24	193/24948 (0.8%)

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	C	0	1
1	D	0	1
All	All	0	4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	458	MET	SD-CE	-5.62	1.65	1.79
1	B	388	HIS	CG-CD2	5.41	1.41	1.35
1	B	286	ALA	CA-CB	-5.17	1.45	1.53
1	B	201	PHE	C-O	-5.05	1.18	1.24
1	A	388	HIS	CG-CD2	5.01	1.41	1.35

All (193) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	513	ARG	N-CA-C	-10.08	96.17	110.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	513	ARG	N-CA-C	-9.92	96.39	110.08
1	A	513	ARG	N-CA-C	-9.70	96.70	110.08
1	C	127	PRO	CA-C-N	9.16	129.09	120.03
1	C	127	PRO	C-N-CA	9.16	129.09	120.03
1	C	513	ARG	N-CA-C	-9.13	97.48	110.08
1	D	137	LYS	N-CA-C	8.43	121.23	111.11
1	A	287	VAL	N-CA-C	8.36	126.74	109.34
1	A	148	TYR	N-CA-C	-8.35	95.93	109.96
1	B	127	PRO	CA-C-N	8.35	128.30	120.03
1	B	127	PRO	C-N-CA	8.35	128.30	120.03
1	A	137	LYS	N-CA-C	8.30	121.07	111.11
1	B	137	LYS	N-CA-C	8.14	120.88	111.11
1	B	208	GLN	N-CA-C	-8.14	103.48	113.41
1	D	127	PRO	CA-C-N	8.12	128.61	119.92
1	D	127	PRO	C-N-CA	8.12	128.61	119.92
1	C	287	VAL	N-CA-C	8.08	126.14	109.34
1	D	388	HIS	CA-CB-CG	-8.05	105.75	113.80
1	A	574	GLY	N-CA-C	-8.03	98.66	112.64
1	B	287	VAL	N-CA-C	8.02	126.02	109.34
1	A	127	PRO	CA-C-N	8.00	127.95	120.03
1	A	127	PRO	C-N-CA	8.00	127.95	120.03
1	C	137	LYS	N-CA-C	7.95	120.65	111.11
1	C	574	GLY	N-CA-C	-7.92	98.86	112.64
1	D	287	VAL	N-CA-C	7.86	125.68	109.34
1	B	233	ILE	N-CA-C	-7.85	105.53	112.12
1	B	388	HIS	CA-CB-CG	-7.71	106.08	113.80
1	D	574	GLY	N-CA-C	-7.60	99.41	112.64
1	B	574	GLY	N-CA-C	-7.54	99.53	112.64
1	A	279	ILE	CA-C-N	7.53	129.25	119.84
1	A	279	ILE	C-N-CA	7.53	129.25	119.84
1	B	279	ILE	CA-C-N	7.48	129.19	119.84
1	B	279	ILE	C-N-CA	7.48	129.19	119.84
1	C	388	HIS	CA-CB-CG	-7.44	106.36	113.80
1	B	258	GLY	N-CA-C	-7.44	98.19	112.02
1	B	148	TYR	N-CA-C	-7.39	97.54	109.96
1	A	388	HIS	CA-CB-CG	-7.38	106.42	113.80
1	B	134	TYR	N-CA-C	7.28	120.65	108.20
1	D	258	GLY	N-CA-C	-7.28	98.48	112.02
1	C	279	ILE	CA-C-N	7.21	128.86	119.84
1	C	279	ILE	C-N-CA	7.21	128.86	119.84
1	C	134	TYR	N-CA-C	7.18	120.49	108.20
1	D	134	TYR	N-CA-C	7.12	120.60	107.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	148	TYR	N-CA-C	-7.11	98.02	109.96
1	D	73	GLU	N-CA-C	-7.07	100.02	110.48
1	D	279	ILE	CA-C-N	7.05	128.65	119.84
1	D	279	ILE	C-N-CA	7.05	128.65	119.84
1	D	208	GLN	N-CA-C	-6.95	102.79	112.45
1	A	208	GLN	N-CA-C	-6.90	103.86	112.90
1	B	73	GLU	N-CA-C	-6.78	100.45	110.48
1	C	148	TYR	N-CA-C	-6.77	98.58	109.96
1	C	258	GLY	N-CA-C	-6.76	99.44	112.02
1	C	73	GLU	N-CA-C	-6.75	100.49	110.48
1	A	258	GLY	N-CA-C	-6.74	99.48	112.02
1	B	482	THR	N-CA-C	6.66	118.54	111.28
1	A	233	ILE	N-CA-C	-6.55	106.62	112.12
1	D	328	LEU	N-CA-C	-6.53	104.09	111.07
1	B	180	LYS	N-CA-C	6.52	118.39	111.28
1	B	77	ARG	N-CA-C	-6.52	103.86	110.97
1	C	208	GLN	N-CA-C	-6.43	104.47	112.90
1	A	134	TYR	N-CA-C	6.38	120.20	107.62
1	A	73	GLU	N-CA-C	-6.33	101.11	110.48
1	B	524	GLU	N-CA-C	6.33	118.98	111.71
1	B	418	GLY	N-CA-C	6.29	128.09	113.18
1	C	41	CYS	N-CA-C	6.29	118.53	108.41
1	D	180	LYS	N-CA-C	6.28	118.64	111.11
1	D	418	GLY	N-CA-C	6.26	128.02	113.18
1	A	180	LYS	N-CA-C	6.24	118.08	111.28
1	C	180	LYS	N-CA-C	6.23	118.07	111.28
1	D	403	SER	N-CA-C	-6.23	102.14	110.55
1	C	418	GLY	N-CA-C	6.21	127.90	113.18
1	A	403	SER	N-CA-C	-6.19	102.19	110.55
1	B	41	CYS	N-CA-C	6.18	118.36	108.41
1	D	142	PHE	N-CA-C	6.07	117.98	111.36
1	A	328	LEU	N-CA-C	-6.06	104.76	111.36
1	A	418	GLY	N-CA-C	6.04	127.49	113.18
1	A	138	SER	N-CA-C	6.01	123.60	110.80
1	C	185	ARG	N-CA-C	-5.96	102.46	111.34
1	B	185	ARG	N-CA-C	-5.96	102.46	111.34
1	B	328	LEU	N-CA-C	-5.94	104.72	111.07
1	A	281	GLU	N-CA-C	-5.93	98.16	110.80
1	A	183	LEU	N-CA-C	5.93	117.95	110.24
1	B	465	GLU	N-CA-C	-5.92	104.75	111.14
1	C	328	LEU	N-CA-C	-5.87	104.79	111.07
1	A	524	GLU	N-CA-C	5.85	118.44	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	233	ILE	N-CA-C	-5.84	106.54	111.56
1	A	422	PHE	N-CA-C	-5.82	98.41	110.80
1	C	370	GLN	N-CA-C	-5.81	99.81	109.46
1	D	226	HIS	N-CA-C	5.81	119.44	111.54
1	C	320	HIS	CA-C-N	5.79	125.64	119.28
1	C	320	HIS	C-N-CA	5.79	125.64	119.28
1	B	388	HIS	ND1-CE1-NE2	5.78	114.18	108.40
1	D	52	PHE	N-CA-C	5.78	118.05	111.11
1	A	478	PHE	N-CA-C	-5.77	105.07	111.71
1	C	63	GLY	N-CA-C	-5.75	107.24	115.64
1	D	465	GLU	N-CA-C	-5.75	104.92	111.07
1	B	370	GLN	N-CA-C	-5.75	100.62	109.76
1	D	550	PHE	N-CA-C	5.69	120.22	112.88
1	D	63	GLY	N-CA-C	-5.69	107.16	115.63
1	C	422	PHE	N-CA-C	-5.68	98.71	110.80
1	B	138	SER	N-CA-C	5.66	122.86	110.80
1	B	503	LEU	N-CA-C	5.66	117.13	111.07
1	D	422	PHE	N-CA-C	-5.65	98.76	110.80
1	D	346	GLU	N-CA-C	5.65	119.89	112.89
1	C	233	ILE	N-CA-C	-5.64	106.71	111.56
1	B	446	ALA	N-CA-C	-5.63	105.22	111.36
1	D	185	ARG	N-CA-C	-5.61	102.49	111.02
1	B	478	PHE	N-CA-C	-5.61	105.17	111.28
1	A	426	PHE	N-CA-C	5.61	118.16	111.71
1	D	138	SER	N-CA-C	5.60	122.73	110.80
1	A	283	LEU	N-CA-C	-5.58	106.15	114.64
1	A	41	CYS	N-CA-C	5.58	117.40	108.41
1	D	306	LEU	N-CA-C	-5.58	105.20	111.28
1	A	77	ARG	N-CA-C	-5.57	104.90	110.97
1	C	215	LYS	N-CA-C	-5.57	105.29	111.36
1	C	201	PHE	N-CA-C	-5.55	105.31	111.36
1	B	142	PHE	N-CA-C	5.55	117.41	111.36
1	A	346	GLU	N-CA-C	5.55	119.89	113.18
1	D	283	LEU	N-CA-C	-5.54	105.50	114.09
1	A	185	ARG	N-CA-C	-5.54	103.09	111.34
1	A	276	PRO	CA-C-N	5.53	126.76	119.84
1	A	276	PRO	C-N-CA	5.53	126.76	119.84
1	D	482	THR	N-CA-C	5.51	118.05	111.71
1	C	465	GLU	N-CA-C	-5.51	105.17	111.07
1	D	183	LEU	N-CA-C	5.51	117.40	110.24
1	A	482	THR	N-CA-C	5.50	118.04	111.71
1	B	346	GLU	N-CA-C	5.49	119.70	112.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	283	LEU	N-CA-C	-5.49	105.58	114.09
1	C	102	ILE	N-CA-C	-5.48	105.38	110.53
1	D	321	PRO	N-CA-C	-5.48	106.72	113.84
1	C	478	PHE	N-CA-C	-5.48	105.31	111.28
1	A	370	GLN	N-CA-C	-5.46	101.08	109.76
1	D	446	ALA	N-CA-C	-5.46	105.41	111.36
1	D	388	HIS	CB-CG-CD2	-5.45	124.12	131.20
1	C	226	HIS	N-CA-C	5.43	118.93	111.54
1	B	422	PHE	N-CA-C	-5.43	99.24	110.80
1	B	388	HIS	ND1-CG-CD2	5.42	111.52	106.10
1	D	289	GLN	N-CA-C	-5.40	100.43	109.24
1	D	41	CYS	N-CA-C	5.39	117.08	108.41
1	C	197	MET	N-CA-C	-5.38	105.42	111.28
1	C	446	ALA	N-CA-C	-5.38	105.42	111.28
1	B	320	HIS	CA-C-N	5.38	125.04	119.56
1	B	320	HIS	C-N-CA	5.38	125.04	119.56
1	B	63	GLY	N-CA-C	-5.36	107.81	115.64
1	A	102	ILE	N-CA-C	-5.36	105.49	110.53
1	C	281	GLU	N-CA-C	-5.35	99.39	110.80
1	D	478	PHE	N-CA-C	-5.35	105.55	111.71
1	B	306	LEU	N-CA-C	-5.35	105.44	111.28
1	D	524	GLU	N-CA-C	5.35	117.87	111.71
1	C	388	HIS	CB-CG-CD2	-5.33	124.27	131.20
1	A	446	ALA	N-CA-C	-5.32	105.56	111.36
1	C	183	LEU	N-CA-C	5.32	117.16	110.24
1	C	546	LYS	CA-C-N	5.28	126.44	119.84
1	C	546	LYS	C-N-CA	5.28	126.44	119.84
1	C	138	SER	N-CA-C	5.28	122.04	110.80
1	A	63	GLY	N-CA-C	-5.27	107.95	115.64
1	C	346	GLU	N-CA-C	5.26	119.55	113.18
1	B	426	PHE	N-CA-C	5.26	117.75	111.71
1	A	257	ILE	N-CA-C	-5.25	100.81	108.17
1	D	105(A)	ILE	CA-C-N	5.21	126.35	119.84
1	D	105(A)	ILE	C-N-CA	5.21	126.35	119.84
1	D	215	LYS	N-CA-C	-5.21	105.28	111.69
1	B	281	GLU	N-CA-C	-5.21	99.71	110.80
1	B	197	MET	N-CA-C	-5.20	105.61	111.28
1	B	388	HIS	CB-CG-CD2	-5.20	124.44	131.20
1	C	276	PRO	CA-C-N	5.20	126.34	119.84
1	C	276	PRO	C-N-CA	5.20	126.34	119.84
1	A	135	GLY	N-CA-C	-5.18	108.53	115.32
1	C	423	VAL	N-CA-C	-5.17	105.67	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	503	LEU	N-CA-C	5.17	116.60	111.07
1	D	320	HIS	CA-C-N	5.16	124.82	119.56
1	D	320	HIS	C-N-CA	5.16	124.82	119.56
1	C	503	LEU	N-CA-C	5.16	116.59	111.07
1	A	388	HIS	ND1-CG-CD2	5.14	111.24	106.10
1	B	283	LEU	N-CA-C	-5.11	106.87	114.64
1	C	306	LEU	N-CA-C	-5.11	105.60	111.07
1	C	550	PHE	N-CA-C	5.11	119.47	112.88
1	D	201	PHE	N-CA-C	-5.11	105.63	111.14
1	B	321	PRO	N-CA-C	-5.10	107.22	113.84
1	D	197	MET	N-CA-C	-5.09	105.73	111.28
1	C	377	ILE	N-CA-CB	5.09	116.57	110.31
1	D	188	ILE	N-CA-C	-5.09	102.07	107.77
1	A	179	GLU	N-CA-C	5.08	116.63	111.14
1	C	403	SER	N-CA-C	-5.08	103.69	110.55
1	D	40	PRO	N-CA-C	5.08	120.44	113.84
1	D	370	GLN	N-CA-C	-5.08	101.03	109.46
1	A	142	PHE	N-CA-C	5.07	116.62	111.14
1	A	404	PHE	N-CA-C	-5.04	104.87	111.02
1	B	286	ALA	N-CA-C	5.03	117.72	111.24
1	C	426	PHE	N-CA-C	5.03	117.49	111.71
1	D	560	ASN	N-CA-C	5.02	119.16	113.19
1	A	388	HIS	CG-CD2-NE2	-5.01	102.19	107.20
1	B	69	CYS	N-CA-C	-5.01	104.48	111.39

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	147	TYR	Sidechain
1	A	466	TYR	Sidechain
1	C	147	TYR	Sidechain
1	D	147	TYR	Sidechain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4473	0	4375	293	0
1	B	4473	0	4375	289	0
1	C	4473	0	4375	291	0
1	D	4473	0	4375	296	0
2	A	42	0	39	1	0
2	B	42	0	39	1	0
2	C	42	0	39	1	0
2	D	42	0	39	0	0
3	A	43	0	30	4	0
3	B	43	0	30	5	0
3	C	43	0	30	5	0
3	D	43	0	30	5	0
4	A	25	0	15	8	0
4	B	25	0	15	9	0
4	C	25	0	15	8	0
4	D	25	0	15	9	0
All	All	18332	0	17836	1155	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 32.

All (1155) 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:477:SER:HB2	1:B:479:GLU:HG2	1.37	1.07
1:C:477:SER:HB2	1:C:479:GLU:HG2	1.41	1.01
1:B:183:LEU:HD22	1:B:442:ILE:HD13	1.44	0.98
1:D:574:GLY:O	1:D:576:PRO:HD3	1.63	0.97
1:B:67:GLU:HB3	2:B:661:NAG:H82	1.44	0.97
1:A:477:SER:HB2	1:A:479:GLU:HG2	1.45	0.96
1:C:183:LEU:HD22	1:C:442:ILE:HD13	1.48	0.95
1:A:574:GLY:O	1:A:576:PRO:HD3	1.66	0.94
1:D:183:LEU:HD22	1:D:442:ILE:HD13	1.48	0.94
1:B:574:GLY:O	1:B:576:PRO:HD3	1.67	0.93
1:D:477:SER:HB2	1:D:479:GLU:HG2	1.50	0.92
1:C:574:GLY:O	1:C:576:PRO:HD3	1.69	0.91
1:A:183:LEU:HD22	1:A:442:ILE:HD13	1.52	0.91
1:A:579:SER:HB2	1:D:267:LYS:HZ2	1.39	0.88
1:B:430:ILE:HG23	1:B:582:VAL:HG11	1.56	0.88
1:C:430:ILE:HG23	1:C:582:VAL:HG11	1.54	0.87
1:A:430:ILE:HG23	1:A:582:VAL:HG11	1.55	0.87
1:A:458:MET:HE2	1:A:460:TYR:HE1	1.37	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:458:MET:HE2	1:B:460:TYR:HE1	1.40	0.86
1:A:156:ALA:HB3	1:A:159:CYS:SG	2.15	0.86
1:D:430:ILE:HG23	1:D:582:VAL:HG11	1.57	0.86
1:C:280:PRO:HG2	1:C:283:LEU:HD12	1.55	0.84
3:A:682:HEM:HBC2	3:A:682:HEM:HHD	1.59	0.84
1:C:458:MET:HE2	1:C:460:TYR:HE1	1.43	0.84
3:D:682:HEM:HBC2	3:D:682:HEM:HHD	1.60	0.84
1:A:579:SER:HB2	1:D:267:LYS:NZ	1.92	0.83
1:A:197:MET:HE2	1:A:578:THR:HG21	1.59	0.82
1:A:280:PRO:HG2	1:A:283:LEU:HD12	1.61	0.82
1:C:156:ALA:HB3	1:C:159:CYS:SG	2.19	0.82
1:C:104:ASN:HB3	1:C:358:LYS:HE3	1.63	0.81
1:D:458:MET:HE2	1:D:460:TYR:HE1	1.46	0.80
1:A:280:PRO:CG	1:A:283:LEU:HD12	2.11	0.80
1:D:472:LEU:HD21	1:D:524:GLU:HG3	1.62	0.80
1:C:553:GLU:HG3	1:C:557:LYS:NZ	1.96	0.79
1:B:120:ARG:HH12	4:B:701:IMN:C18	1.96	0.79
1:D:553:GLU:HG3	1:D:557:LYS:NZ	1.96	0.79
1:C:75:LEU:HA	1:C:78:ILE:HD12	1.65	0.78
1:A:75:LEU:HA	1:A:78:ILE:HD12	1.65	0.78
1:B:291:VAL:HG22	1:B:294:LEU:HD12	1.66	0.78
1:A:553:GLU:HG3	1:A:557:LYS:NZ	1.99	0.78
3:B:682:HEM:HBC2	3:B:682:HEM:HHD	1.64	0.77
1:C:133:HIS:HD2	1:C:147:TYR:OH	1.67	0.77
1:D:454:GLN:HA	1:D:457:GLU:HG2	1.66	0.77
1:C:477:SER:CB	1:C:479:GLU:HG2	2.15	0.77
1:A:104:ASN:HB3	1:A:358:LYS:HE3	1.67	0.77
1:C:352:LEU:O	4:C:701:IMN:H63	1.85	0.77
1:C:472:LEU:HD21	1:C:524:GLU:HG3	1.66	0.76
1:D:108:LEU:O	1:D:112:ILE:HG12	1.86	0.76
1:D:156:ALA:HB3	1:D:159:CYS:SG	2.25	0.76
1:A:382:ASN:HD21	3:A:682:HEM:HAD2	1.51	0.76
1:C:197:MET:HE2	1:C:578:THR:HG21	1.68	0.75
1:D:197:MET:HE2	1:D:578:THR:HG21	1.68	0.75
1:B:192:GLN:OE1	1:B:517:ILE:HG22	1.86	0.75
1:B:230:LEU:HD13	1:B:233:ILE:HD12	1.67	0.74
1:B:280:PRO:HG3	1:B:283:LEU:HD12	1.66	0.74
1:C:454:GLN:HA	1:C:457:GLU:HG2	1.69	0.74
1:D:291:VAL:HG22	1:D:294:LEU:HD12	1.68	0.74
1:D:75:LEU:HA	1:D:78:ILE:HD12	1.67	0.74
1:D:218:PRO:HB2	1:D:458:MET:SD	2.27	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:LEU:HA	1:B:78:ILE:HD12	1.69	0.74
1:B:472:LEU:HD21	1:B:524:GLU:HG3	1.68	0.74
1:A:108:LEU:O	1:A:112:ILE:HG12	1.88	0.74
1:A:349:VAL:HG22	4:A:701:IMN:O1	1.88	0.73
1:B:281:GLU:O	1:B:283:LEU:N	2.20	0.73
1:A:577:PHE:HE2	1:D:267:LYS:HD3	1.51	0.73
1:B:280:PRO:CG	1:B:283:LEU:HD12	2.18	0.73
1:C:192:GLN:OE1	1:C:517:ILE:HG22	1.88	0.73
1:C:242:HIS:CE1	1:C:245:ARG:HH21	2.07	0.73
3:C:682:HEM:HBC2	3:C:682:HEM:HHD	1.71	0.72
1:A:291:VAL:HG22	1:A:294:LEU:HD12	1.71	0.72
1:C:123:LEU:O	1:C:469:ARG:NH2	2.22	0.72
1:D:352:LEU:O	4:D:701:IMN:H63	1.89	0.72
1:A:109:ARG:HG3	1:A:357:PHE:CE1	2.25	0.72
1:D:75:LEU:HG	1:D:78:ILE:HD12	1.70	0.72
1:C:280:PRO:CG	1:C:283:LEU:HD12	2.19	0.72
1:B:156:ALA:HB3	1:B:159:CYS:SG	2.30	0.71
1:A:472:LEU:HD21	1:A:524:GLU:HG3	1.70	0.71
1:D:230:LEU:HD13	1:D:233:ILE:HD12	1.72	0.71
1:A:211:LYS:NZ	1:A:236:GLU:HG3	2.04	0.71
1:C:75:LEU:HG	1:C:78:ILE:HD12	1.70	0.71
1:B:266:VAL:HG13	1:B:271:VAL:O	1.91	0.71
1:B:120:ARG:NH1	4:B:701:IMN:C18	2.53	0.71
1:B:197:MET:HE2	1:B:578:THR:HG21	1.72	0.71
1:A:192:GLN:OE1	1:A:517:ILE:HG22	1.91	0.71
1:A:281:GLU:O	1:A:283:LEU:N	2.24	0.71
1:C:291:VAL:HG22	1:C:294:LEU:HD12	1.71	0.71
1:A:123:LEU:O	1:A:469:ARG:NH2	2.24	0.70
1:D:526:GLY:HA3	4:D:701:IMN:C14	2.22	0.70
1:B:108:LEU:O	1:B:112:ILE:HG12	1.91	0.70
1:B:211:LYS:NZ	1:B:236:GLU:HG3	2.07	0.70
1:B:477:SER:CB	1:B:479:GLU:HG2	2.19	0.70
1:C:230:LEU:HD13	1:C:233:ILE:HD12	1.71	0.70
1:C:211:LYS:NZ	1:C:236:GLU:HG3	2.07	0.70
1:D:192:GLN:OE1	1:D:517:ILE:HG22	1.92	0.70
1:A:275:TYR:CD1	1:A:279:ILE:HD12	2.27	0.70
1:B:454:GLN:HA	1:B:457:GLU:HG2	1.71	0.69
1:C:189:PRO:CB	1:C:430:ILE:HD12	2.21	0.69
1:D:420:THR:HG22	1:D:576:PRO:HG3	1.74	0.69
1:D:109:ARG:HG3	1:D:357:PHE:CE1	2.27	0.69
1:C:150:ARG:NH2	1:C:458:MET:O	2.26	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:ASN:HB3	1:B:358:LYS:HE3	1.73	0.69
1:B:442:ILE:O	1:B:445:GLN:HG2	1.93	0.69
1:C:430:ILE:CG2	1:C:582:VAL:HG11	2.22	0.69
1:B:75:LEU:HG	1:B:78:ILE:HD12	1.73	0.69
1:D:281:GLU:O	1:D:283:LEU:N	2.26	0.68
1:A:230:LEU:HD13	1:A:233:ILE:HD12	1.76	0.68
1:C:281:GLU:O	1:C:283:LEU:N	2.26	0.68
1:C:75:LEU:HA	1:C:78:ILE:CD1	2.24	0.68
1:D:384:LEU:HB3	4:D:701:IMN:CL	2.30	0.68
1:B:352:LEU:O	4:B:701:IMN:H63	1.93	0.68
1:D:83:LYS:HA	1:D:83:LYS:HE3	1.76	0.68
1:C:252:LEU:HD22	1:C:309:HIS:CD2	2.29	0.68
1:C:108:LEU:O	1:C:112:ILE:HG12	1.93	0.68
1:C:211:LYS:HZ2	1:C:236:GLU:HG3	1.59	0.68
1:B:150:ARG:NH2	1:B:458:MET:O	2.27	0.68
1:A:191:PRO:HD2	1:A:433:ARG:HG3	1.75	0.68
1:A:577:PHE:CE2	1:D:267:LYS:HD3	2.29	0.68
1:B:104:ASN:HB3	1:B:358:LYS:CE	2.23	0.68
1:D:430:ILE:CG2	1:D:582:VAL:HG11	2.23	0.67
1:B:218:PRO:HB2	1:B:458:MET:SD	2.33	0.67
1:B:133:HIS:HD2	1:B:147:TYR:OH	1.75	0.67
1:B:184:ARG:HA	1:B:438:ARG:O	1.94	0.67
1:B:553:GLU:HG3	1:B:557:LYS:NZ	2.10	0.67
1:C:553:GLU:HG3	1:C:557:LYS:HZ3	1.56	0.67
1:A:252:LEU:HD22	1:A:309:HIS:CD2	2.30	0.67
1:D:211:LYS:NZ	1:D:236:GLU:HG3	2.09	0.67
1:D:266:VAL:HG13	1:D:271:VAL:O	1.94	0.67
1:D:189:PRO:CB	1:D:430:ILE:HD12	2.24	0.67
1:A:189:PRO:HB3	1:A:430:ILE:HD12	1.76	0.67
1:C:478:PHE:CE2	1:C:495:TYR:HB2	2.30	0.67
1:A:294:LEU:HA	1:A:409:TYR:HB3	1.77	0.67
1:D:553:GLU:HG3	1:D:557:LYS:HZ3	1.57	0.67
1:A:75:LEU:HG	1:A:78:ILE:HD12	1.77	0.67
1:C:109:ARG:HG3	1:C:357:PHE:CE1	2.30	0.67
1:B:128:PRO:HG3	1:B:376:ARG:HH11	1.59	0.66
1:C:189:PRO:HB3	1:C:430:ILE:HD12	1.75	0.66
1:A:430:ILE:CG2	1:A:582:VAL:HG11	2.25	0.66
1:A:33:ALA:HB3	1:A:158:ASP:OD2	1.95	0.66
1:B:478:PHE:CE2	1:B:495:TYR:HB2	2.30	0.66
1:D:104:ASN:HB3	1:D:358:LYS:HE3	1.77	0.66
1:D:280:PRO:CG	1:D:283:LEU:HD12	2.26	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:294:LEU:HA	1:D:409:TYR:HB3	1.77	0.66
1:D:189:PRO:HB3	1:D:430:ILE:HD12	1.77	0.66
1:B:123:LEU:O	1:B:469:ARG:NH2	2.29	0.65
1:C:353:SER:HA	4:C:701:IMN:O	1.96	0.65
1:C:33:ALA:HB3	1:C:158:ASP:OD2	1.97	0.65
1:D:252:LEU:HD22	1:D:309:HIS:CD2	2.31	0.65
1:B:182:LEU:O	1:B:438:ARG:HA	1.97	0.65
1:A:487:MET:O	1:A:490:GLU:HB3	1.97	0.65
1:B:430:ILE:CG2	1:B:582:VAL:HG11	2.27	0.65
1:D:485:LYS:HA	1:D:488:ALA:HB3	1.78	0.65
1:A:477:SER:CB	1:A:479:GLU:HG2	2.22	0.65
1:C:294:LEU:HA	1:C:409:TYR:HB3	1.78	0.65
1:D:442:ILE:O	1:D:445:GLN:HG2	1.97	0.65
1:D:123:LEU:O	1:D:469:ARG:NH2	2.29	0.64
1:B:137:LYS:HB2	1:B:137:LYS:NZ	2.12	0.64
1:B:252:LEU:HD22	1:B:309:HIS:CD2	2.32	0.64
1:B:485:LYS:HA	1:B:488:ALA:HB3	1.79	0.64
1:A:479:GLU:CD	1:A:479:GLU:H	2.04	0.64
1:B:75:LEU:HA	1:B:78:ILE:CD1	2.26	0.64
1:D:150:ARG:NH2	1:D:458:MET:O	2.31	0.64
1:A:479:GLU:HB3	1:A:485:LYS:NZ	2.13	0.64
1:B:148:TYR:HD1	1:B:377:ILE:HG12	1.63	0.64
1:A:319:GLU:HG3	1:A:554:VAL:HG11	1.79	0.64
1:D:478:PHE:CE2	1:D:495:TYR:HB2	2.33	0.64
1:A:485:LYS:HA	1:A:488:ALA:HB3	1.78	0.64
1:B:281:GLU:C	1:B:283:LEU:H	2.05	0.64
1:A:150:ARG:NH2	1:A:458:MET:O	2.31	0.64
1:D:485:LYS:HD2	1:D:489:ALA:HB2	1.80	0.63
1:B:213:ASP:OD1	1:B:215:LYS:HE2	1.98	0.63
1:B:482:THR:HG22	1:B:509:VAL:HG12	1.79	0.63
1:C:303:THR:HG22	1:C:307:ARG:HH11	1.63	0.63
1:A:189:PRO:CB	1:A:430:ILE:HD12	2.28	0.63
1:A:424:GLU:HA	1:A:428:ARG:HH21	1.62	0.63
1:C:582:VAL:O	1:C:582:VAL:HG13	1.98	0.63
1:D:353:SER:HA	4:D:701:IMN:O	1.98	0.63
1:A:303:THR:HG22	1:A:307:ARG:HH11	1.64	0.63
1:D:120:ARG:HH12	4:D:701:IMN:C18	2.10	0.63
1:C:148:TYR:HD1	1:C:377:ILE:HG12	1.64	0.63
1:A:478:PHE:CE2	1:A:495:TYR:HB2	2.34	0.63
1:B:109:ARG:HG3	1:B:357:PHE:CE1	2.33	0.63
1:D:303:THR:HG22	1:D:307:ARG:HH11	1.63	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:75:LEU:HA	1:D:78:ILE:CD1	2.28	0.62
1:A:109:ARG:HG3	1:A:357:PHE:HE1	1.64	0.62
1:A:104:ASN:HB3	1:A:358:LYS:CE	2.28	0.62
1:B:189:PRO:HB3	1:B:430:ILE:HD12	1.80	0.62
1:A:122:TYR:O	1:A:122:TYR:CD1	2.53	0.62
1:A:75:LEU:HA	1:A:78:ILE:CD1	2.29	0.62
1:B:294:LEU:HA	1:B:409:TYR:HB3	1.82	0.62
1:B:256:VAL:HA	1:B:260:GLU:O	1.99	0.62
1:C:137:LYS:HB2	1:C:137:LYS:NZ	2.15	0.62
1:C:420:THR:HG22	1:C:576:PRO:HG3	1.82	0.62
1:C:479:GLU:HB3	1:C:485:LYS:CE	2.29	0.62
1:D:128:PRO:HG3	1:D:376:ARG:HH11	1.64	0.62
1:A:442:ILE:O	1:A:445:GLN:HG2	2.00	0.62
1:B:145:LEU:HD23	1:B:376:ARG:NH2	2.15	0.62
1:C:242:HIS:HE1	1:C:245:ARG:HH21	1.44	0.62
1:C:330:GLN:HB3	1:D:138:SER:HB2	1.81	0.62
1:D:104:ASN:HB3	1:D:358:LYS:CE	2.29	0.62
1:C:439:ASN:O	1:C:441:PRO:HD3	2.00	0.61
1:D:477:SER:CB	1:D:479:GLU:HG2	2.27	0.61
1:A:525:LEU:N	1:A:525:LEU:HD23	2.15	0.61
1:B:463:LEU:HD22	1:B:506:ALA:CB	2.30	0.61
1:C:122:TYR:CD1	1:C:122:TYR:O	2.54	0.61
1:A:509:VAL:HG12	1:A:509:VAL:O	2.00	0.61
1:A:565:GLN:HG3	1:D:268:ASP:OD2	2.00	0.61
1:B:189:PRO:CB	1:B:430:ILE:HD12	2.31	0.61
1:B:319:GLU:HG3	1:B:554:VAL:HG11	1.83	0.61
1:A:190:ASP:OD1	1:A:517:ILE:HB	2.01	0.61
1:A:388:HIS:CE1	1:A:447:VAL:HG11	2.35	0.61
1:D:137:LYS:NZ	1:D:137:LYS:HB2	2.15	0.61
1:A:137:LYS:HB2	1:A:137:LYS:NZ	2.16	0.61
1:B:424:GLU:HA	1:B:428:ARG:HH21	1.65	0.61
1:B:303:THR:HG22	1:B:307:ARG:HH11	1.65	0.61
1:B:582:VAL:O	1:B:582:VAL:HG13	1.99	0.61
1:C:479:GLU:CD	1:C:479:GLU:H	2.08	0.61
1:D:191:PRO:HD2	1:D:433:ARG:HG3	1.82	0.61
1:A:428:ARG:HA	1:A:582:VAL:HG23	1.83	0.61
1:B:525:LEU:N	1:B:525:LEU:HD23	2.15	0.61
1:C:479:GLU:HB3	1:C:485:LYS:NZ	2.16	0.61
1:D:275:TYR:CD1	1:D:279:ILE:HD12	2.36	0.61
1:C:213:ASP:OD1	1:C:215:LYS:HE2	2.01	0.60
1:C:353:SER:HA	4:C:701:IMN:C3	2.30	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:420:THR:HG22	1:A:576:PRO:HG3	1.82	0.60
1:C:320:HIS:HE1	1:C:551:GLY:O	1.84	0.60
1:D:33:ALA:HB3	1:D:158:ASP:OD2	2.02	0.60
1:D:582:VAL:O	1:D:582:VAL:HG13	2.01	0.60
1:A:148:TYR:HD1	1:A:377:ILE:HG12	1.66	0.60
1:A:506:ALA:O	1:A:510:GLU:HB2	2.02	0.60
1:D:506:ALA:O	1:D:510:GLU:HB2	2.01	0.60
1:C:104:ASN:HB3	1:C:358:LYS:CE	2.30	0.60
1:C:218:PRO:HB2	1:C:458:MET:SD	2.41	0.60
1:C:230:LEU:HG	1:C:337:ILE:HG12	1.84	0.60
1:C:557:LYS:N	1:C:557:LYS:HE3	2.16	0.60
1:D:405:LYS:H	1:D:405:LYS:HD2	1.67	0.60
1:B:33:ALA:HB3	1:B:158:ASP:OD2	2.01	0.60
1:B:420:THR:HG22	1:B:576:PRO:HG3	1.84	0.60
1:C:388:HIS:CE1	1:C:447:VAL:HG11	2.37	0.60
1:B:510:GLU:O	1:B:512:PRO:HD3	2.02	0.60
1:A:330:GLN:HB3	1:B:138:SER:HB2	1.84	0.59
1:C:491:LEU:HD11	1:C:509:VAL:HG11	1.84	0.59
1:A:34:ASN:HB3	1:A:37:CYS:SG	2.42	0.59
1:A:275:TYR:HD1	1:A:279:ILE:HD12	1.65	0.59
1:D:234:TYR:CE1	1:D:252:LEU:HD21	2.37	0.59
1:A:479:GLU:HB3	1:A:485:LYS:CE	2.32	0.59
1:A:454:GLN:HA	1:A:457:GLU:HG2	1.83	0.59
1:A:582:VAL:O	1:A:582:VAL:HG13	2.00	0.59
1:C:349:VAL:HG22	4:C:701:IMN:O1	2.03	0.59
1:A:230:LEU:HG	1:A:337:ILE:HG12	1.85	0.59
1:A:491:LEU:HD11	1:A:509:VAL:HG11	1.83	0.59
1:B:342:LYS:HG2	1:B:562:ALA:HB3	1.84	0.59
1:B:482:THR:OG1	1:B:488:ALA:HB2	2.03	0.59
1:C:256:VAL:HA	1:C:260:GLU:O	2.02	0.59
1:C:266:VAL:HG13	1:C:271:VAL:O	2.03	0.59
1:D:391:LEU:HD21	3:D:682:HEM:HAB	1.85	0.59
1:A:276:PRO:HG2	1:A:409:TYR:CD1	2.38	0.59
1:A:553:GLU:HG3	1:A:557:LYS:HZ3	1.66	0.59
1:B:185:ARG:HE	1:B:438:ARG:HD3	1.68	0.59
1:C:124:ILE:HD11	1:C:528:PRO:C	2.28	0.59
1:C:485:LYS:HA	1:C:488:ALA:HB3	1.84	0.59
1:A:463:LEU:HD22	1:A:506:ALA:CB	2.33	0.59
1:B:88:THR:O	1:B:91:TYR:HB3	2.03	0.59
1:C:504:TYR:HB3	1:C:505:PRO:HD3	1.85	0.59
1:D:230:LEU:HG	1:D:337:ILE:HG12	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:557:LYS:HE3	1:D:557:LYS:N	2.17	0.59
1:A:241:GLN:HG3	1:A:242:HIS:N	2.18	0.59
1:A:281:GLU:C	1:A:283:LEU:H	2.10	0.59
1:B:479:GLU:HA	1:B:488:ALA:HB1	1.85	0.59
1:D:463:LEU:HD22	1:D:506:ALA:CB	2.33	0.59
1:D:353:SER:HA	4:D:701:IMN:C6	2.33	0.58
1:D:424:GLU:HA	1:D:428:ARG:HH21	1.67	0.58
1:A:94:THR:HG23	1:A:354:GLY:O	2.03	0.58
1:A:211:LYS:HZ2	1:A:236:GLU:HG3	1.67	0.58
1:A:218:PRO:HB2	1:A:458:MET:SD	2.44	0.58
1:C:253:LYS:HD3	1:C:269:THR:HG22	1.85	0.58
1:D:145:LEU:HD23	1:D:376:ARG:NH2	2.17	0.58
1:B:491:LEU:HD11	1:B:509:VAL:HG11	1.85	0.58
1:C:120:ARG:HH12	4:C:701:IMN:C18	2.16	0.58
1:D:182:LEU:O	1:D:438:ARG:HA	2.02	0.58
1:B:405:LYS:H	1:B:405:LYS:HD2	1.68	0.58
1:A:134:TYR:HD2	1:A:136:TYR:CE1	2.22	0.58
1:A:554:VAL:HG13	1:A:555:GLY:H	1.69	0.58
1:B:349:VAL:HG22	4:B:701:IMN:O1	2.04	0.58
1:C:191:PRO:HG3	1:C:433:ARG:CZ	2.33	0.58
1:A:128:PRO:HG3	1:A:376:ARG:HH11	1.68	0.58
1:C:190:ASP:OD1	1:C:517:ILE:HB	2.03	0.58
1:B:83:LYS:HE3	1:B:83:LYS:HA	1.85	0.58
1:C:83:LYS:HE3	1:C:83:LYS:HA	1.86	0.58
1:C:442:ILE:O	1:C:445:GLN:HG2	2.03	0.58
1:B:230:LEU:HG	1:B:337:ILE:HG12	1.85	0.58
1:B:541:SER:HB2	1:B:543:GLN:OE1	2.04	0.58
1:A:485:LYS:HD2	1:A:489:ALA:HB2	1.86	0.58
1:A:145:LEU:HD23	1:A:376:ARG:NH2	2.19	0.57
1:A:306:LEU:HD23	1:A:306:LEU:C	2.29	0.57
1:B:478:PHE:CZ	1:B:495:TYR:HB2	2.39	0.57
1:C:420:THR:CG2	1:C:576:PRO:HG3	2.34	0.57
1:D:213:ASP:OD1	1:D:215:LYS:HE2	2.05	0.57
1:D:525:LEU:HD23	1:D:525:LEU:N	2.19	0.57
1:B:506:ALA:O	1:B:510:GLU:HB2	2.05	0.57
1:C:182:LEU:O	1:C:438:ARG:HA	2.04	0.57
1:D:183:LEU:O	1:D:438:ARG:HB3	2.05	0.57
1:B:428:ARG:HA	1:B:582:VAL:HG23	1.87	0.57
1:C:281:GLU:C	1:C:283:LEU:H	2.12	0.57
1:A:120:ARG:NH1	4:A:701:IMN:C18	2.68	0.57
1:A:178:LEU:O	1:A:182:LEU:HB2	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:253:LYS:HD3	1:D:269:THR:HG22	1.85	0.57
1:D:280:PRO:HG3	1:D:283:LEU:HD12	1.85	0.57
1:D:420:THR:CG2	1:D:576:PRO:HG3	2.33	0.57
1:C:478:PHE:CZ	1:C:495:TYR:HB2	2.40	0.57
1:D:124:ILE:HD11	1:D:528:PRO:C	2.29	0.57
1:D:479:GLU:HA	1:D:488:ALA:HB1	1.87	0.57
1:A:120:ARG:HH12	4:A:701:IMN:C18	2.16	0.57
1:A:114:LYS:HD3	1:A:369:GLN:NE2	2.19	0.57
1:C:506:ALA:O	1:C:510:GLU:HB2	2.04	0.57
1:D:198:PHE:HB2	1:D:580:PHE:HB3	1.86	0.57
1:A:198:PHE:HB2	1:A:580:PHE:HB3	1.87	0.56
1:C:191:PRO:HD2	1:C:433:ARG:HG3	1.86	0.56
1:C:543:GLN:NE2	1:D:127:PRO:O	2.38	0.56
1:B:179:GLU:HA	1:B:183:LEU:HD12	1.87	0.56
1:B:242:HIS:CE1	1:B:245:ARG:HH21	2.23	0.56
1:C:184:ARG:HA	1:C:438:ARG:O	2.04	0.56
3:D:682:HEM:HHD	3:D:682:HEM:CBC	2.33	0.56
1:A:482:THR:HG22	1:A:509:VAL:HG12	1.86	0.56
1:B:276:PRO:HG2	1:B:409:TYR:CD1	2.40	0.56
1:D:120:ARG:NH1	4:D:701:IMN:C18	2.68	0.56
1:D:405:LYS:H	1:D:405:LYS:CD	2.18	0.56
1:A:462:SER:OG	1:A:465:GLU:HG2	2.05	0.56
1:D:256:VAL:HA	1:D:260:GLU:O	2.06	0.56
1:B:241:GLN:HG3	1:B:242:HIS:N	2.20	0.56
1:B:291:VAL:CG2	1:B:294:LEU:HD12	2.35	0.56
1:C:114:LYS:HD3	1:C:369:GLN:NE2	2.20	0.56
1:D:485:LYS:O	1:D:485:LYS:HG3	2.05	0.56
1:B:72:PRO:HB2	1:B:76:THR:HB	1.88	0.56
1:B:128:PRO:HG3	1:B:376:ARG:NH1	2.19	0.56
1:B:462:SER:OG	1:B:465:GLU:HG2	2.06	0.56
1:C:128:PRO:HG3	1:C:376:ARG:HH11	1.69	0.56
1:C:412:SER:O	1:C:416:GLU:N	2.38	0.56
1:D:482:THR:HG22	1:D:509:VAL:HG12	1.87	0.56
1:A:553:GLU:HG3	1:A:557:LYS:HZ1	1.68	0.56
1:C:241:GLN:O	1:C:245:ARG:HG3	2.06	0.56
1:C:475:TYR:CD1	1:C:480:GLU:HG2	2.41	0.56
1:B:124:ILE:HD11	1:B:528:PRO:C	2.30	0.56
1:B:134:TYR:HD2	1:B:136:TYR:CE1	2.23	0.56
1:B:464:ASN:HD21	1:B:475:TYR:H	1.52	0.56
1:B:553:GLU:HG3	1:B:557:LYS:HZ1	1.70	0.56
1:C:105:ASN:O	1:C:106:PRO:HD3	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:148:TYR:HD1	1:D:377:ILE:HG12	1.71	0.56
1:D:479:GLU:CD	1:D:479:GLU:H	2.14	0.56
1:D:527:ALA:HA	4:D:701:IMN:H162	1.87	0.56
1:A:124:ILE:HD11	1:A:528:PRO:C	2.30	0.56
1:A:216:ARG:HB3	1:A:220:PHE:CD2	2.41	0.56
1:A:482:THR:OG1	1:A:488:ALA:HB2	2.06	0.56
1:B:211:LYS:HZ2	1:B:236:GLU:HG3	1.70	0.56
1:B:487:MET:O	1:B:490:GLU:HB3	2.05	0.56
1:C:568:ILE:HG23	1:C:572:VAL:HG21	1.86	0.56
1:C:482:THR:HG22	1:C:509:VAL:HG12	1.86	0.56
1:A:458:MET:HE2	1:A:460:TYR:CE1	2.29	0.55
1:B:306:LEU:C	1:B:306:LEU:HD23	2.30	0.55
1:C:276:PRO:HG2	1:C:409:TYR:CD1	2.41	0.55
1:C:134:TYR:HD2	1:C:136:TYR:CE1	2.23	0.55
1:C:477:SER:HB2	1:C:479:GLU:CG	2.24	0.55
1:B:122:TYR:O	1:B:122:TYR:CD1	2.59	0.55
1:C:145:LEU:HD23	1:C:376:ARG:NE	2.21	0.55
1:C:463:LEU:HD22	1:C:506:ALA:CB	2.35	0.55
1:A:242:HIS:CE1	1:A:245:ARG:HH21	2.24	0.55
1:D:281:GLU:C	1:D:283:LEU:H	2.13	0.55
1:A:88:THR:O	1:A:92:ILE:HG13	2.07	0.55
1:A:478:PHE:CZ	1:A:495:TYR:HB2	2.42	0.55
1:A:504:TYR:HB3	1:A:505:PRO:HD3	1.87	0.55
1:C:479:GLU:HB3	1:C:485:LYS:HE3	1.89	0.55
1:D:72:PRO:HB2	1:D:76:THR:HB	1.87	0.55
1:D:83:LYS:HA	1:D:83:LYS:CE	2.36	0.55
1:D:241:GLN:HG3	1:D:242:HIS:N	2.19	0.55
1:B:568:ILE:HG23	1:B:572:VAL:HG21	1.87	0.55
1:C:88:THR:O	1:C:92:ILE:HG13	2.06	0.55
1:C:184:ARG:NH1	1:C:441:PRO:HG3	2.22	0.55
1:C:464:ASN:HD21	1:C:475:TYR:H	1.54	0.55
1:D:260:GLU:HB2	1:D:262:TYR:HE2	1.72	0.55
1:A:292:PHE:O	1:A:299:MET:HE2	2.07	0.55
1:A:479:GLU:HA	1:A:488:ALA:HB1	1.89	0.55
1:A:554:VAL:HG13	1:A:555:GLY:N	2.21	0.55
1:C:72:PRO:HB2	1:C:76:THR:HB	1.89	0.55
1:D:242:HIS:CE1	1:D:245:ARG:HH21	2.25	0.55
1:D:388:HIS:CE1	1:D:447:VAL:HG11	2.41	0.55
1:A:211:LYS:HZ1	1:A:236:GLU:HG3	1.72	0.55
1:B:198:PHE:CD1	1:B:198:PHE:C	2.85	0.55
1:B:388:HIS:CE1	1:B:447:VAL:HG11	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:287:VAL:HG11	1:C:299:MET:SD	2.47	0.55
1:C:405:LYS:CD	1:C:405:LYS:H	2.20	0.55
1:D:74:PHE:HB2	1:D:77:ARG:NH1	2.21	0.55
1:D:122:TYR:O	1:D:122:TYR:CD1	2.59	0.55
1:D:160:PRO:HG2	1:D:164:GLY:O	2.07	0.55
1:D:382:ASN:HD21	3:D:682:HEM:HAD2	1.72	0.55
1:D:464:ASN:HD21	1:D:475:TYR:H	1.55	0.55
1:D:482:THR:OG1	1:D:488:ALA:HB2	2.06	0.55
1:A:479:GLU:HB3	1:A:485:LYS:HE3	1.89	0.55
1:B:198:PHE:HB2	1:B:580:PHE:HB3	1.89	0.55
1:B:420:THR:CG2	1:B:576:PRO:HG3	2.37	0.55
1:B:424:GLU:O	1:B:428:ARG:NE	2.40	0.55
1:C:198:PHE:HB2	1:C:580:PHE:HB3	1.89	0.55
1:D:34:ASN:HB3	1:D:37:CYS:SG	2.47	0.55
1:D:306:LEU:C	1:D:306:LEU:HD23	2.32	0.55
1:A:128:PRO:HG3	1:A:376:ARG:NH1	2.22	0.55
1:C:482:THR:OG1	1:C:488:ALA:HB2	2.07	0.55
1:A:475:TYR:CD1	1:A:480:GLU:HG2	2.42	0.54
1:A:120:ARG:HG2	1:A:531:LEU:HD12	1.90	0.54
1:B:509:VAL:HG12	1:B:509:VAL:O	2.06	0.54
1:A:424:GLU:O	1:A:428:ARG:NE	2.41	0.54
1:B:405:LYS:H	1:B:405:LYS:CD	2.19	0.54
1:B:463:LEU:HD22	1:B:506:ALA:HB1	1.89	0.54
1:A:184:ARG:NH1	1:A:441:PRO:HG3	2.23	0.54
1:A:568:ILE:O	1:A:572:VAL:HG23	2.07	0.54
1:B:320:HIS:HE1	1:B:551:GLY:O	1.91	0.54
1:C:306:LEU:HD23	1:C:306:LEU:C	2.33	0.54
1:C:495:TYR:CE2	1:C:501:MET:SD	3.00	0.54
1:A:266:VAL:HG13	1:A:271:VAL:O	2.08	0.54
1:B:423:VAL:HG13	1:B:578:THR:CG2	2.38	0.54
1:D:380:GLU:HG2	1:D:466:TYR:CE1	2.43	0.54
1:D:275:TYR:HD1	1:D:279:ILE:HD12	1.71	0.54
1:D:342:LYS:HG2	1:D:562:ALA:HB3	1.89	0.54
1:A:565:GLN:HE22	1:A:577:PHE:HA	1.73	0.54
1:B:137:LYS:HB2	1:B:137:LYS:HZ3	1.72	0.54
1:B:458:MET:HE2	1:B:460:TYR:CE1	2.32	0.54
1:C:148:TYR:CD1	1:C:377:ILE:HG12	2.43	0.54
1:C:234:TYR:CE1	1:C:252:LEU:HD21	2.43	0.54
1:D:216:ARG:HB3	1:D:220:PHE:CD2	2.43	0.54
1:B:353:SER:HA	4:B:701:IMN:C6	2.37	0.54
1:D:105:ASN:O	1:D:106:PRO:HD3	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:128:PRO:HG3	1:D:376:ARG:NH1	2.22	0.54
3:A:682:HEM:HHD	3:A:682:HEM:CBC	2.37	0.54
1:C:273:MET:SD	1:C:290:GLU:HA	2.47	0.54
1:C:463:LEU:HD22	1:C:506:ALA:HB1	1.90	0.54
1:D:46:GLU:O	1:D:57:CYS:HA	2.07	0.54
1:D:320:HIS:HE1	1:D:551:GLY:O	1.91	0.54
1:D:478:PHE:CZ	1:D:495:TYR:HB2	2.43	0.54
1:A:420:THR:CG2	1:A:576:PRO:HG3	2.37	0.54
1:B:475:TYR:CD1	1:B:480:GLU:HG2	2.42	0.54
1:C:261:VAL:O	1:C:307:ARG:NH1	2.41	0.54
1:D:191:PRO:HG3	1:D:433:ARG:CZ	2.38	0.54
1:D:491:LEU:HD11	1:D:509:VAL:HG11	1.88	0.54
1:A:439:ASN:O	1:A:441:PRO:HD3	2.08	0.53
1:B:292:PHE:O	1:B:299:MET:HE2	2.08	0.53
1:B:557:LYS:HE3	1:B:557:LYS:N	2.23	0.53
1:C:67:GLU:HB3	2:C:661:NAG:H82	1.90	0.53
1:D:276:PRO:HG2	1:D:409:TYR:CD1	2.43	0.53
1:D:315:ILE:HG21	1:D:558:ILE:HD11	1.89	0.53
1:A:256:VAL:HA	1:A:260:GLU:O	2.08	0.53
1:C:315:ILE:HG21	1:C:558:ILE:HD11	1.90	0.53
1:A:83:LYS:HE3	1:A:83:LYS:HA	1.88	0.53
1:A:213:ASP:HB2	1:A:222:ARG:HG3	1.91	0.53
1:A:405:LYS:CD	1:A:405:LYS:H	2.22	0.53
1:D:245:ARG:HD3	1:D:329:PHE:CE1	2.44	0.53
1:B:94:THR:HG23	1:B:354:GLY:O	2.09	0.53
1:B:150:ARG:NH1	1:B:150:ARG:HG2	2.23	0.53
1:B:269:THR:OG1	1:B:271:VAL:HG13	2.09	0.53
1:B:209:PHE:HB2	1:B:377:ILE:HG21	1.91	0.53
1:C:216:ARG:HB3	1:C:220:PHE:CD2	2.44	0.53
1:A:72:PRO:HB2	1:A:76:THR:HB	1.90	0.53
1:B:260:GLU:HB2	1:B:262:TYR:HE2	1.73	0.53
1:A:423:VAL:HG13	1:A:578:THR:CG2	2.37	0.53
1:B:216:ARG:HB3	1:B:220:PHE:CD2	2.44	0.53
1:C:213:ASP:HB2	1:C:222:ARG:HG3	1.90	0.53
1:C:241:GLN:HG3	1:C:242:HIS:N	2.24	0.53
1:C:479:GLU:HA	1:C:488:ALA:HB1	1.90	0.53
1:D:145:LEU:HD23	1:D:376:ARG:NE	2.24	0.53
1:A:70:THR:OG1	2:A:661:NAG:H81	2.09	0.53
1:A:245:ARG:HH22	1:A:326:GLU:HG2	1.73	0.53
1:C:275:TYR:CD1	1:C:279:ILE:HD12	2.44	0.53
1:A:568:ILE:HG23	1:A:572:VAL:HG21	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:539:ILE:HA	1:B:544:TYR:HB3	1.91	0.53
1:C:105(A):ILE:HG21	1:C:108:LEU:HD12	1.90	0.53
1:C:210:PHE:CE1	1:C:382:ASN:HA	2.44	0.53
1:C:89:VAL:HG12	1:C:93:LEU:CD2	2.39	0.52
1:C:213:ASP:OD1	1:C:215:LYS:HB2	2.09	0.52
1:C:292:PHE:O	1:C:299:MET:HE2	2.09	0.52
1:D:88:THR:O	1:D:91:TYR:HB3	2.09	0.52
1:A:191:PRO:HG3	1:A:433:ARG:CZ	2.39	0.52
1:A:210:PHE:CE1	1:A:382:ASN:HA	2.44	0.52
1:A:260:GLU:HB2	1:A:262:TYR:HE2	1.74	0.52
1:A:303:THR:CG2	1:A:307:ARG:HH11	2.22	0.52
1:B:211:LYS:HZ1	1:B:236:GLU:HG3	1.74	0.52
1:B:353:SER:HA	4:B:701:IMN:O	2.10	0.52
1:C:195:ASN:HB3	1:C:582:VAL:HB	1.92	0.52
1:C:291:VAL:CG2	1:C:294:LEU:HD12	2.39	0.52
1:D:149:THR:HG21	1:D:376:ARG:HB3	1.91	0.52
1:B:148:TYR:CD1	1:B:377:ILE:HG12	2.42	0.52
1:C:128:PRO:HG3	1:C:376:ARG:NH1	2.25	0.52
1:C:275:TYR:HD1	1:C:279:ILE:HD12	1.73	0.52
1:D:85:THR:O	1:D:89:VAL:HG23	2.09	0.52
1:D:190:ASP:OD1	1:D:517:ILE:HB	2.10	0.52
1:D:509:VAL:HG12	1:D:509:VAL:O	2.10	0.52
1:B:64:PHE:CE2	1:B:72:PRO:HB3	2.44	0.52
1:B:124:ILE:HD12	1:B:124:ILE:N	2.25	0.52
1:D:50:THR:CG2	1:D:56:LYS:HB2	2.39	0.52
1:D:145:LEU:HD23	1:D:376:ARG:CZ	2.40	0.52
1:A:320:HIS:HE1	1:A:551:GLY:O	1.92	0.52
1:B:74:PHE:HB2	1:B:77:ARG:NH1	2.25	0.52
1:B:105:ASN:O	1:B:106:PRO:HD3	2.10	0.52
1:C:94:THR:HG23	1:C:354:GLY:O	2.08	0.52
1:D:211:LYS:HZ2	1:D:236:GLU:HG3	1.75	0.52
1:D:291:VAL:CG2	1:D:294:LEU:HD12	2.37	0.52
1:B:412:SER:O	1:B:416:GLU:N	2.43	0.52
1:C:260:GLU:HB2	1:C:262:TYR:HE2	1.74	0.52
1:C:543:GLN:HE22	1:D:127:PRO:HD2	1.74	0.52
1:B:145:LEU:HD23	1:B:376:ARG:HH21	1.75	0.52
1:D:198:PHE:CD1	1:D:198:PHE:C	2.88	0.52
1:D:352:LEU:HD21	1:D:387:TRP:CH2	2.45	0.52
1:A:182:LEU:O	1:A:438:ARG:HA	2.10	0.52
1:A:242:HIS:HE1	1:A:245:ARG:HH21	1.56	0.52
1:A:281:GLU:HA	1:A:284:GLN:HG3	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:245:ARG:HD3	1:B:329:PHE:CE1	2.44	0.52
1:C:414:LEU:HA	1:C:422:PHE:CE2	2.45	0.52
1:D:211:LYS:HZ1	1:D:236:GLU:HG3	1.75	0.52
1:D:412:SER:O	1:D:416:GLU:N	2.42	0.52
1:C:335:ILE:HA	1:C:559:ILE:HD11	1.92	0.52
1:D:133:HIS:HD2	1:D:147:TYR:OH	1.93	0.52
1:D:150:ARG:NH1	1:D:150:ARG:HG2	2.24	0.52
1:D:303:THR:CG2	1:D:307:ARG:HH11	2.22	0.52
1:C:391:LEU:HD21	3:C:682:HEM:HAB	1.91	0.51
1:D:91:TYR:O	1:D:95:HIS:HD2	1.93	0.51
1:A:145:LEU:HD23	1:A:376:ARG:CZ	2.40	0.51
1:A:280:PRO:HG3	1:A:283:LEU:HD12	1.88	0.51
1:C:189:PRO:HB2	1:C:430:ILE:HD12	1.90	0.51
1:D:50:THR:HG21	1:D:56:LYS:HB2	1.91	0.51
1:A:464:ASN:HD21	1:A:475:TYR:H	1.56	0.51
1:B:398:GLU:HG2	1:B:421:GLN:NE2	2.24	0.51
1:C:198:PHE:C	1:C:198:PHE:CD1	2.89	0.51
1:C:303:THR:CG2	1:C:307:ARG:HH11	2.22	0.51
1:D:245:ARG:NH2	1:D:247:PHE:HE1	2.09	0.51
1:A:477:SER:HB2	1:A:479:GLU:CG	2.30	0.51
1:A:578:THR:O	1:D:267:LYS:NZ	2.43	0.51
1:C:462:SER:OG	1:C:465:GLU:HG2	2.11	0.51
1:D:109:ARG:HG3	1:D:357:PHE:HE1	1.71	0.51
1:C:85:THR:O	1:C:89:VAL:HG23	2.11	0.51
1:C:179:GLU:HA	1:C:183:LEU:HD12	1.93	0.51
1:D:88:THR:O	1:D:92:ILE:HG13	2.10	0.51
1:D:504:TYR:HB3	1:D:505:PRO:HD3	1.93	0.51
1:A:148:TYR:CD1	1:A:377:ILE:HG12	2.44	0.51
1:A:320:HIS:HD2	1:B:49:SER:O	1.94	0.51
1:A:372:GLN:HB3	1:A:374:GLN:HG2	1.92	0.51
1:B:315:ILE:HG21	1:B:558:ILE:HD11	1.91	0.51
1:C:424:GLU:HA	1:C:428:ARG:HH21	1.75	0.51
1:B:190:ASP:OD1	1:B:517:ILE:HB	2.10	0.51
1:D:134:TYR:HD2	1:D:136:TYR:CE1	2.29	0.51
1:D:280:PRO:HG2	1:D:283:LEU:HD12	1.92	0.51
1:D:439:ASN:O	1:D:441:PRO:HD3	2.11	0.51
1:A:89:VAL:HG12	1:A:93:LEU:CD2	2.40	0.51
1:B:479:GLU:CD	1:B:479:GLU:H	2.18	0.51
1:B:479:GLU:HB3	1:B:485:LYS:CE	2.41	0.51
1:B:89:VAL:HG12	1:B:93:LEU:CD2	2.41	0.51
1:B:372:GLN:HB3	1:B:374:GLN:HG2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:362:ASP:O	1:C:365:LEU:HG	2.11	0.51
1:A:276:PRO:O	1:A:279:ILE:HG13	2.10	0.51
1:A:412:SER:O	1:A:416:GLU:N	2.41	0.51
1:B:145:LEU:HD23	1:B:376:ARG:NE	2.25	0.51
1:B:234:TYR:CE1	1:B:252:LEU:HD21	2.46	0.51
1:B:353:SER:HA	4:B:701:IMN:H63	1.93	0.51
1:B:452:ILE:O	1:B:456:ARG:HG3	2.11	0.51
1:C:372:GLN:HB3	1:C:374:GLN:HG2	1.93	0.51
1:C:382:ASN:HD21	3:C:682:HEM:HAD2	1.76	0.51
1:D:64:PHE:CE2	1:D:72:PRO:HB3	2.46	0.51
1:B:479:GLU:HB3	1:B:485:LYS:NZ	2.27	0.50
1:C:48:MET:O	1:C:55:TYR:HA	2.11	0.50
1:A:322:GLU:HG2	1:B:52:PHE:HA	1.93	0.50
1:A:510:GLU:O	1:A:512:PRO:HD3	2.11	0.50
1:C:428:ARG:HA	1:C:582:VAL:HG23	1.93	0.50
1:C:458:MET:HE2	1:C:460:TYR:CE1	2.34	0.50
1:C:539:ILE:HA	1:C:544:TYR:HB3	1.93	0.50
1:A:91:TYR:O	1:A:95:HIS:HD2	1.94	0.50
1:B:372:GLN:OE1	1:B:374:GLN:NE2	2.44	0.50
1:B:565:GLN:HE22	1:B:577:PHE:HA	1.75	0.50
1:C:380:GLU:HG2	1:C:466:TYR:CE1	2.47	0.50
1:D:210:PHE:CE1	1:D:382:ASN:HA	2.46	0.50
1:D:541:SER:HB2	1:D:543:GLN:OE1	2.10	0.50
1:A:97:LYS:HB2	1:A:356:HIS:CE1	2.46	0.50
1:C:396:ASN:ND2	1:C:401:GLU:HB3	2.27	0.50
1:D:131:ASN:ND2	1:D:147:TYR:CD2	2.79	0.50
1:A:245:ARG:HD3	1:A:329:PHE:CD1	2.46	0.50
1:A:315:ILE:HG21	1:A:558:ILE:HD11	1.94	0.50
1:B:46:GLU:O	1:B:57:CYS:HA	2.11	0.50
1:B:439:ASN:O	1:B:441:PRO:HD3	2.12	0.50
1:C:405:LYS:H	1:C:405:LYS:HD2	1.76	0.50
1:C:64:PHE:CE2	1:C:72:PRO:HB3	2.47	0.50
1:C:294:LEU:HD23	1:C:409:TYR:HB3	1.93	0.50
1:D:276:PRO:HB2	1:D:278:HIS:CE1	2.47	0.50
1:B:145:LEU:HD23	1:B:376:ARG:CZ	2.41	0.50
1:B:382:ASN:HD21	3:B:682:HEM:HAD2	1.77	0.50
1:D:554:VAL:HG13	1:D:555:GLY:N	2.26	0.50
1:B:261:VAL:O	1:B:307:ARG:NH1	2.45	0.50
1:D:179:GLU:HA	1:D:183:LEU:HD12	1.93	0.50
1:A:198:PHE:CD1	1:A:198:PHE:C	2.90	0.50
1:B:160:PRO:HG2	1:B:164:GLY:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:113:MET:O	1:C:116:VAL:HG22	2.11	0.50
1:C:510:GLU:O	1:C:512:PRO:HD3	2.12	0.50
1:D:178:LEU:O	1:D:182:LEU:HB2	2.12	0.50
1:D:396:ASN:ND2	1:D:401:GLU:HB3	2.27	0.50
1:D:539:ILE:HA	1:D:544:TYR:HB3	1.94	0.50
1:A:191:PRO:CD	1:A:433:ARG:HG3	2.42	0.49
1:C:42:GLN:HB2	1:C:68:ASN:O	2.12	0.49
1:C:352:LEU:HD21	1:C:387:TRP:CH2	2.47	0.49
1:D:462:SER:OG	1:D:465:GLU:HG2	2.12	0.49
1:A:145:LEU:HD23	1:A:376:ARG:NE	2.27	0.49
1:A:234:TYR:CE1	1:A:252:LEU:HD21	2.47	0.49
1:A:564:ILE:HG12	1:A:580:PHE:CE1	2.47	0.49
1:B:388:HIS:N	1:B:389:PRO:CD	2.76	0.49
1:C:160:PRO:HG2	1:C:164:GLY:O	2.12	0.49
1:C:553:GLU:HG3	1:C:557:LYS:HZ1	1.76	0.49
1:D:423:VAL:HG13	1:D:578:THR:CG2	2.41	0.49
1:D:428:ARG:HA	1:D:582:VAL:HG23	1.93	0.49
1:A:352:LEU:HD21	1:A:387:TRP:CH2	2.47	0.49
1:B:472:LEU:HD11	1:B:524:GLU:HB2	1.95	0.49
1:A:150:ARG:NH1	1:A:150:ARG:HG2	2.25	0.49
1:B:191:PRO:HG3	1:B:433:ARG:CZ	2.42	0.49
1:C:180:LYS:HD3	1:C:490:GLU:CD	2.37	0.49
1:C:525:LEU:N	1:C:525:LEU:HD23	2.27	0.49
1:D:261:VAL:O	1:D:307:ARG:NH1	2.45	0.49
1:D:452:ILE:O	1:D:456:ARG:HG3	2.12	0.49
1:A:261:VAL:O	1:A:307:ARG:NH1	2.45	0.49
1:D:510:GLU:O	1:D:512:PRO:HD3	2.11	0.49
1:D:553:GLU:HG3	1:D:557:LYS:HZ1	1.76	0.49
1:A:213:ASP:OD1	1:A:215:LYS:HE2	2.11	0.49
1:A:291:VAL:CG2	1:A:294:LEU:HD12	2.42	0.49
1:B:380:GLU:O	1:B:384:LEU:HG	2.12	0.49
1:C:150:ARG:NH1	1:C:150:ARG:HG2	2.26	0.49
1:B:353:SER:HA	4:B:701:IMN:C3	2.42	0.49
1:C:423:VAL:HG13	1:C:578:THR:CG2	2.42	0.49
1:D:149:THR:O	1:D:378:ALA:HA	2.13	0.49
1:D:319:GLU:HG3	1:D:554:VAL:HG11	1.95	0.49
1:B:88:THR:O	1:B:92:ILE:HG13	2.13	0.49
1:B:281:GLU:C	1:B:283:LEU:N	2.69	0.49
1:C:145:LEU:HD23	1:C:376:ARG:NH2	2.27	0.49
1:C:160:PRO:HG2	1:C:165:VAL:HA	1.95	0.49
1:D:522:MET:O	4:D:701:IMN:H14	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:453:ASP:O	1:A:456:ARG:HB2	2.13	0.49
1:B:253:LYS:HD3	1:B:269:THR:HG22	1.94	0.49
1:C:109:ARG:HG3	1:C:357:PHE:HE1	1.75	0.49
1:C:462:SER:HB2	1:C:499:ASP:O	2.13	0.49
1:C:565:GLN:HE22	1:C:577:PHE:HA	1.78	0.49
1:D:155:VAL:H	1:D:459:LYS:HB3	1.78	0.49
1:A:83:LYS:HA	1:A:83:LYS:CE	2.43	0.49
1:D:94:THR:HG23	1:D:354:GLY:O	2.13	0.49
1:D:241:GLN:O	1:D:245:ARG:HG3	2.12	0.49
1:D:434:VAL:H	1:D:517:ILE:HG13	1.77	0.49
1:D:472:LEU:CD2	1:D:524:GLU:HG3	2.38	0.49
1:A:575:CYS:O	1:A:576:PRO:C	2.54	0.48
1:C:544:TYR:OH	1:D:142:PHE:HB2	2.13	0.48
1:A:384:LEU:HD23	1:A:507:LEU:HD11	1.94	0.48
1:B:210:PHE:CE1	1:B:382:ASN:HA	2.48	0.48
1:C:353:SER:HA	4:C:701:IMN:C6	2.44	0.48
1:C:509:VAL:HG12	1:C:509:VAL:O	2.14	0.48
1:D:184:ARG:NH1	1:D:441:PRO:HG3	2.28	0.48
1:C:145:LEU:HD23	1:C:376:ARG:CZ	2.42	0.48
1:D:89:VAL:HG12	1:D:93:LEU:CD2	2.43	0.48
1:B:275:TYR:HD1	1:B:279:ILE:HD12	1.77	0.48
1:D:113:MET:O	1:D:116:VAL:HG22	2.14	0.48
1:B:554:VAL:HG13	1:B:555:GLY:N	2.29	0.48
1:C:65:TYR:HE1	1:C:73:GLU:HA	1.78	0.48
1:B:303:THR:CG2	1:B:307:ARG:HH11	2.26	0.48
1:B:380:GLU:HG2	1:B:466:TYR:CE1	2.49	0.48
1:D:380:GLU:O	1:D:384:LEU:HG	2.14	0.48
1:D:414:LEU:HA	1:D:422:PHE:CE2	2.48	0.48
1:A:138:SER:HB2	1:B:330:GLN:HB3	1.95	0.48
1:A:184:ARG:HA	1:A:438:ARG:O	2.13	0.48
1:D:114:LYS:HD3	1:D:369:GLN:NE2	2.29	0.48
1:D:116:VAL:O	1:D:120:ARG:HB2	2.14	0.48
1:D:160:PRO:HG2	1:D:165:VAL:HA	1.95	0.48
1:D:287:VAL:HG11	1:D:299:MET:SD	2.53	0.48
1:D:568:ILE:HG23	1:D:572:VAL:HG21	1.95	0.48
1:A:380:GLU:HG2	1:A:466:TYR:CE1	2.49	0.48
1:A:413:ILE:O	1:A:416:GLU:HB2	2.13	0.48
1:B:109:ARG:HG3	1:B:357:PHE:HE1	1.76	0.48
1:C:138:SER:HB2	1:D:330:GLN:HB3	1.95	0.48
1:A:211:LYS:O	1:A:212:THR:C	2.57	0.48
1:A:463:LEU:O	1:A:467:ARG:HG3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:ARG:HG2	1:B:531:LEU:HD12	1.96	0.48
1:B:396:ASN:HA	1:B:401:GLU:HA	1.96	0.48
1:D:244:LEU:HD11	1:D:288:GLY:HA2	1.96	0.48
1:A:105(A):ILE:HG21	1:A:108:LEU:HD12	1.96	0.48
1:A:127:PRO:HD2	1:B:543:GLN:HE22	1.79	0.48
1:A:195:ASN:HB3	1:A:582:VAL:HB	1.95	0.48
1:A:414:LEU:HA	1:A:422:PHE:CE2	2.49	0.48
1:B:150:ARG:HH11	1:B:150:ARG:CG	2.27	0.48
1:C:413:ILE:O	1:C:416:GLU:HB2	2.14	0.48
1:D:475:TYR:CD1	1:D:480:GLU:HG2	2.49	0.48
1:C:410:ASN:OD1	1:C:413:ILE:HG13	2.13	0.47
1:C:485:LYS:HD2	1:C:489:ALA:HB2	1.95	0.47
1:C:575:CYS:O	1:C:576:PRO:C	2.56	0.47
1:A:160:PRO:HG2	1:A:165:VAL:HA	1.95	0.47
1:B:113:MET:HE2	1:B:360:LYS:O	2.13	0.47
1:B:124:ILE:HD12	1:B:124:ILE:H	1.78	0.47
1:D:137:LYS:O	1:D:138:SER:O	2.31	0.47
1:A:380:GLU:O	1:A:384:LEU:HG	2.14	0.47
1:B:114:LYS:HE2	1:B:365:LEU:O	2.14	0.47
1:B:137:LYS:O	1:B:138:SER:O	2.31	0.47
1:C:178:LEU:O	1:C:182:LEU:HB2	2.15	0.47
3:C:682:HEM:HHD	3:C:682:HEM:CBC	2.41	0.47
1:D:372:GLN:HB3	1:D:374:GLN:HG2	1.96	0.47
1:A:160:PRO:HG2	1:A:164:GLY:O	2.14	0.47
1:A:179:GLU:HA	1:A:183:LEU:HD12	1.95	0.47
1:B:113:MET:HG2	1:B:360:LYS:HB3	1.96	0.47
1:B:160:PRO:HG2	1:B:165:VAL:HA	1.96	0.47
1:B:244:LEU:HD11	1:B:288:GLY:HA2	1.97	0.47
1:B:459:LYS:HE2	1:B:459:LYS:HA	1.96	0.47
1:B:564:ILE:HG12	1:B:580:PHE:CE1	2.50	0.47
1:C:384:LEU:HD23	1:C:507:LEU:HD11	1.95	0.47
1:D:113:MET:HE3	1:D:117:LEU:HD11	1.96	0.47
1:D:189:PRO:HB2	1:D:430:ILE:HD12	1.96	0.47
1:D:389:PRO:HB2	1:D:434:VAL:HA	1.97	0.47
1:A:463:LEU:HD22	1:A:506:ALA:HB1	1.96	0.47
1:A:543:GLN:NE2	1:B:127:PRO:O	2.48	0.47
1:B:427:THR:OG1	1:B:578:THR:HG22	2.15	0.47
1:C:453:ASP:O	1:C:456:ARG:HB2	2.13	0.47
1:C:485:LYS:O	1:C:485:LYS:HG3	2.13	0.47
1:D:478:PHE:CE1	1:D:498:ILE:HA	2.49	0.47
1:D:554:VAL:HG13	1:D:555:GLY:H	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:244:LEU:HD11	1:C:288:GLY:HA2	1.97	0.47
1:D:120:ARG:HG2	1:D:531:LEU:HD12	1.97	0.47
1:D:213:ASP:HB2	1:D:222:ARG:HG3	1.96	0.47
1:D:433:ARG:NH2	1:D:512:PRO:HB2	2.29	0.47
1:A:244:LEU:HD11	1:A:288:GLY:HA2	1.95	0.47
1:A:342:LYS:HG2	1:A:562:ALA:HB3	1.97	0.47
1:A:464:ASN:OD1	1:A:499:ASP:HA	2.14	0.47
1:A:526:GLY:HA3	4:A:701:IMN:C14	2.44	0.47
1:B:195:ASN:HB3	1:B:582:VAL:HB	1.96	0.47
3:B:682:HEM:HHD	3:B:682:HEM:CBC	2.39	0.47
1:C:424:GLU:O	1:C:428:ARG:NE	2.48	0.47
1:C:568:ILE:O	1:C:572:VAL:HG23	2.14	0.47
1:D:180:LYS:HD2	1:D:494:LEU:HD11	1.96	0.47
1:D:209:PHE:HB2	1:D:377:ILE:HG21	1.97	0.47
1:D:472:LEU:HD21	1:D:524:GLU:CG	2.40	0.47
1:A:105:ASN:O	1:A:106:PRO:HD3	2.15	0.47
1:B:385:TYR:CZ	4:B:701:IMN:H12	2.50	0.47
1:B:472:LEU:HD21	1:B:524:GLU:CG	2.42	0.47
1:D:124:ILE:N	1:D:124:ILE:HD12	2.29	0.47
1:D:464:ASN:OD1	1:D:499:ASP:HA	2.14	0.47
1:B:116:VAL:O	1:B:120:ARG:HB2	2.15	0.47
1:C:113:MET:HE2	1:C:360:LYS:O	2.15	0.47
1:C:564:ILE:HG12	1:C:580:PHE:CE1	2.50	0.47
1:D:464:ASN:ND2	1:D:475:TYR:H	2.13	0.47
1:A:410:ASN:OD1	1:A:413:ILE:HG13	2.14	0.47
1:B:56:LYS:HD3	1:B:57:CYS:N	2.30	0.47
1:C:133:HIS:CD2	1:C:147:TYR:CE2	3.03	0.47
1:C:254:TYR:HA	1:C:264:PRO:HD3	1.96	0.47
1:C:391:LEU:HB2	1:C:441:PRO:HG2	1.97	0.47
1:D:114:LYS:NZ	1:D:368:ASN:HD21	2.13	0.47
1:A:65:TYR:O	1:A:71:THR:HG23	2.15	0.46
1:A:389:PRO:HB2	1:A:434:VAL:HA	1.96	0.46
1:A:405:LYS:H	1:A:405:LYS:HD2	1.79	0.46
1:B:479:GLU:HB3	1:B:485:LYS:HE3	1.96	0.46
1:C:120:ARG:NH1	4:C:701:IMN:C18	2.78	0.46
1:D:126:SER:HB2	1:D:532:LYS:NZ	2.30	0.46
1:D:211:LYS:O	1:D:212:THR:C	2.58	0.46
1:A:46:GLU:O	1:A:57:CYS:HA	2.16	0.46
1:A:396:ASN:HA	1:A:401:GLU:HA	1.96	0.46
1:B:423:VAL:HG13	1:B:578:THR:HG23	1.96	0.46
1:B:477:SER:HB2	1:B:479:GLU:CG	2.26	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:463:LEU:HD22	1:D:506:ALA:HB1	1.97	0.46
1:A:394:THR:HA	1:A:402:TYR:O	2.14	0.46
1:A:423:VAL:HG13	1:A:578:THR:HG23	1.96	0.46
1:B:333:LYS:O	1:B:337:ILE:HG13	2.15	0.46
1:A:539:ILE:HA	1:A:544:TYR:HB3	1.97	0.46
1:B:85:THR:O	1:B:89:VAL:HG23	2.15	0.46
1:C:116:VAL:O	1:C:120:ARG:HB2	2.16	0.46
1:C:211:LYS:O	1:C:212:THR:C	2.57	0.46
1:D:388:HIS:HB3	1:D:444:VAL:HG21	1.98	0.46
1:D:424:GLU:O	1:D:428:ARG:NE	2.49	0.46
1:D:470:PHE:CG	1:D:525:LEU:HD22	2.51	0.46
1:A:76:THR:O	1:A:80:LEU:HG	2.15	0.46
1:A:131:ASN:OD1	1:A:131:ASN:C	2.58	0.46
1:B:245:ARG:HD3	1:B:329:PHE:CD1	2.50	0.46
1:B:504:TYR:HB3	1:B:505:PRO:HD3	1.98	0.46
1:C:122:TYR:CD1	1:C:122:TYR:C	2.93	0.46
1:C:174:SER:OG	1:C:449:LYS:HE2	2.15	0.46
1:C:447:VAL:O	1:C:450:ALA:HB3	2.15	0.46
1:D:279:ILE:HG23	1:D:280:PRO:HD2	1.97	0.46
1:D:345:ILE:HG12	1:D:534:LEU:HD23	1.97	0.46
1:C:120:ARG:HG2	1:C:531:LEU:HD12	1.98	0.46
1:C:286:ALA:O	1:C:287:VAL:HG22	2.16	0.46
1:C:403:SER:O	1:C:404:PHE:C	2.56	0.46
1:C:472:LEU:CD2	1:C:524:GLU:HG3	2.41	0.46
1:A:113:MET:HE3	1:A:117:LEU:HD11	1.96	0.46
1:A:122:TYR:CD1	1:A:122:TYR:C	2.94	0.46
1:A:427:THR:OG1	1:A:578:THR:HG22	2.15	0.46
1:C:113:MET:HE3	1:C:117:LEU:HD11	1.97	0.46
1:C:137:LYS:HB2	1:C:137:LYS:HZ3	1.81	0.46
1:C:276:PRO:HB2	1:C:278:HIS:CE1	2.51	0.46
1:C:472:LEU:HD21	1:C:524:GLU:CG	2.43	0.46
1:D:150:ARG:HH11	1:D:150:ARG:CG	2.28	0.46
1:A:485:LYS:HG3	1:A:485:LYS:O	2.14	0.46
1:C:149:THR:O	1:C:378:ALA:HA	2.16	0.46
1:D:184:ARG:HA	1:D:438:ARG:O	2.16	0.46
1:D:372:GLN:OE1	1:D:374:GLN:NE2	2.48	0.46
1:D:396:ASN:HA	1:D:401:GLU:HA	1.98	0.46
3:D:682:HEM:CMB	3:D:682:HEM:HBB2	2.45	0.46
1:A:51:GLY:C	1:B:322:GLU:HG2	2.41	0.46
1:B:279:ILE:HG23	1:B:280:PRO:HD2	1.98	0.46
1:D:294:LEU:HD23	1:D:409:TYR:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:MET:HE3	1:A:117:LEU:CD1	2.46	0.46
1:B:145:LEU:O	1:B:224:LEU:HD23	2.16	0.46
1:B:495:TYR:CE2	1:B:501:MET:SD	3.08	0.46
1:D:148:TYR:CD1	1:D:377:ILE:HG12	2.51	0.46
1:D:260:GLU:HB2	1:D:262:TYR:CE2	2.49	0.46
1:D:421:GLN:O	1:D:422:PHE:HB2	2.15	0.46
1:D:433:ARG:HH21	1:D:512:PRO:HB2	1.81	0.46
1:D:479:GLU:HB3	1:D:485:LYS:NZ	2.31	0.46
1:A:49:SER:O	1:B:320:HIS:HD2	1.98	0.45
1:B:48:MET:O	1:B:55:TYR:HA	2.17	0.45
1:C:245:ARG:HD3	1:C:329:PHE:CE1	2.51	0.45
1:C:106:PRO:O	1:C:107:PHE:C	2.59	0.45
1:C:396:ASN:HA	1:C:401:GLU:HA	1.98	0.45
1:A:150:ARG:CG	1:A:150:ARG:HH11	2.29	0.45
1:B:379:SER:HB3	1:B:460:TYR:OH	2.17	0.45
1:C:150:ARG:CG	1:C:150:ARG:HH11	2.29	0.45
1:C:281:GLU:C	1:C:283:LEU:N	2.74	0.45
1:A:116:VAL:O	1:A:120:ARG:HB2	2.15	0.45
1:A:403:SER:O	1:A:404:PHE:C	2.60	0.45
1:A:428:ARG:HA	1:A:583:GLN:OE1	2.17	0.45
1:C:83:LYS:HA	1:C:83:LYS:CE	2.46	0.45
1:C:388:HIS:N	1:C:389:PRO:CD	2.80	0.45
1:A:557:LYS:HD3	1:A:557:LYS:HA	1.76	0.45
1:C:452:ILE:O	1:C:456:ARG:HG3	2.17	0.45
1:A:64:PHE:CE2	1:A:72:PRO:HB3	2.51	0.45
1:B:46:GLU:OE2	1:B:137:LYS:HE3	2.17	0.45
1:B:215:LYS:HD3	1:B:215:LYS:N	2.32	0.45
1:B:477:SER:O	1:B:480:GLU:N	2.50	0.45
1:C:74:PHE:HD1	1:C:77:ARG:HH11	1.63	0.45
1:C:132:VAL:HG21	1:C:219:GLY:HA3	1.99	0.45
1:C:379:SER:HB3	1:C:460:TYR:OH	2.16	0.45
1:C:389:PRO:HB2	1:C:434:VAL:HA	1.99	0.45
1:C:464:ASN:HD21	1:C:475:TYR:N	2.15	0.45
1:C:487:MET:O	1:C:490:GLU:HB3	2.16	0.45
1:D:245:ARG:HD3	1:D:329:PHE:CD1	2.52	0.45
1:A:75:LEU:O	1:A:79:LYS:HG3	2.17	0.45
1:A:293:GLY:HA2	1:A:299:MET:HE2	1.98	0.45
1:C:46:GLU:O	1:C:57:CYS:HA	2.17	0.45
1:C:131:ASN:ND2	1:C:147:TYR:CD2	2.85	0.45
1:C:333:LYS:O	1:C:337:ILE:HG13	2.17	0.45
1:D:105(A):ILE:HG21	1:D:108:LEU:HD12	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:122:TYR:CD1	1:D:122:TYR:C	2.95	0.45
1:A:286:ALA:O	1:A:287:VAL:HG22	2.16	0.45
1:B:478:PHE:HE2	1:B:495:TYR:HB2	1.80	0.45
1:C:126:SER:HB2	1:C:532:LYS:NZ	2.32	0.45
1:C:282:ASN:ND2	1:C:283:LEU:HG	2.31	0.45
1:C:554:VAL:HG13	1:C:555:GLY:N	2.32	0.45
1:D:149:THR:CG2	1:D:376:ARG:HB3	2.47	0.45
1:D:458:MET:HE2	1:D:460:TYR:CE1	2.37	0.45
1:D:575:CYS:O	1:D:576:PRO:C	2.55	0.45
1:A:137:LYS:HB2	1:A:137:LYS:HZ3	1.82	0.45
1:A:353:SER:HA	4:A:701:IMN:H63	1.99	0.45
1:C:91:TYR:CE2	1:C:95:HIS:NE2	2.85	0.45
1:D:433:ARG:HH21	1:D:512:PRO:CB	2.30	0.45
1:A:124:ILE:HD12	1:A:124:ILE:H	1.82	0.45
1:C:459:LYS:HE2	1:C:459:LYS:HA	1.99	0.45
1:D:126:SER:HB2	1:D:532:LYS:HZ3	1.82	0.45
1:D:366:LEU:HA	1:D:369:GLN:HG2	1.98	0.45
1:D:447:VAL:O	1:D:450:ALA:HB3	2.16	0.45
1:C:380:GLU:O	1:C:384:LEU:HG	2.16	0.44
1:C:396:ASN:HD22	1:C:401:GLU:HB3	1.81	0.44
1:B:133:HIS:CD2	1:B:147:TYR:CE2	3.05	0.44
1:D:145:LEU:HD23	1:D:376:ARG:HH21	1.79	0.44
1:A:124:ILE:HD12	1:A:124:ILE:N	2.31	0.44
1:A:352:LEU:HA	1:A:352:LEU:HD12	1.79	0.44
1:B:38:SER:O	1:B:39:ASN:C	2.61	0.44
1:B:76:THR:O	1:B:80:LEU:HG	2.17	0.44
1:B:211:LYS:O	1:B:212:THR:C	2.61	0.44
1:D:403:SER:O	1:D:404:PHE:C	2.61	0.44
1:A:74:PHE:HB2	1:A:77:ARG:NH1	2.32	0.44
1:A:345:ILE:HG12	1:A:534:LEU:HD23	2.00	0.44
1:A:353:SER:HA	4:A:701:IMN:C6	2.47	0.44
1:A:379:SER:HB3	1:A:460:TYR:OH	2.18	0.44
1:B:362:ASP:O	1:B:365:LEU:HG	2.17	0.44
1:B:389:PRO:HB2	1:B:434:VAL:HA	1.98	0.44
1:C:183:LEU:O	1:C:438:ARG:HB3	2.17	0.44
1:B:150:ARG:NH1	1:B:150:ARG:CG	2.81	0.44
1:B:193:GLY:O	1:B:582:VAL:HG12	2.18	0.44
1:B:245:ARG:NH2	1:B:247:PHE:HE1	2.15	0.44
1:B:414:LEU:HA	1:B:422:PHE:CE2	2.53	0.44
1:B:472:LEU:CD2	1:B:524:GLU:HG3	2.41	0.44
1:B:568:ILE:O	1:B:572:VAL:HG23	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:577:PHE:CD1	1:B:577:PHE:C	2.96	0.44
1:C:114:LYS:HD3	1:C:369:GLN:HE21	1.81	0.44
1:C:372:GLN:OE1	1:C:374:GLN:NE2	2.50	0.44
1:D:43:ASN:O	1:D:44:ARG:CB	2.66	0.44
1:D:292:PHE:O	1:D:299:MET:HE2	2.16	0.44
1:A:245:ARG:HH22	1:A:326:GLU:CG	2.30	0.44
1:A:388:HIS:N	1:A:389:PRO:CD	2.80	0.44
1:A:434:VAL:H	1:A:517:ILE:HG13	1.83	0.44
1:A:487:MET:HA	1:A:490:GLU:HB3	2.00	0.44
1:B:133:HIS:CD2	1:B:147:TYR:HE2	2.35	0.44
1:B:183:LEU:O	1:B:438:ARG:HB3	2.18	0.44
1:B:275:TYR:CD1	1:B:279:ILE:HD12	2.52	0.44
1:B:293:GLY:HA2	1:B:299:MET:HE2	2.00	0.44
1:C:320:HIS:HB3	1:C:323:TRP:CG	2.52	0.44
1:D:245:ARG:HH22	1:D:326:GLU:HG2	1.82	0.44
1:D:472:LEU:HD11	1:D:524:GLU:HB2	2.00	0.44
1:D:479:GLU:HA	1:D:488:ALA:CB	2.47	0.44
1:A:245:ARG:HD3	1:A:329:PHE:CE1	2.53	0.44
1:B:339:GLU:HG2	1:B:562:ALA:HB2	1.99	0.44
1:B:410:ASN:OD1	1:B:413:ILE:HG13	2.17	0.44
1:C:34:ASN:HB3	1:C:37:CYS:SG	2.58	0.44
1:C:427:THR:OG1	1:C:578:THR:HG22	2.17	0.44
1:D:137:LYS:HB2	1:D:137:LYS:HZ3	1.81	0.44
1:D:213:ASP:OD1	1:D:215:LYS:HB2	2.18	0.44
1:D:231:ASN:OD1	1:D:231:ASN:C	2.61	0.44
1:D:394:THR:HA	1:D:402:TYR:O	2.17	0.44
1:A:281:GLU:C	1:A:283:LEU:N	2.74	0.44
1:A:398:GLU:O	1:A:399:ASP:HB2	2.18	0.44
1:B:43:ASN:O	1:B:44:ARG:CB	2.66	0.44
1:B:113:MET:HE3	1:B:117:LEU:HD11	2.00	0.44
1:B:345:ILE:HG12	1:B:534:LEU:HD23	2.00	0.44
1:B:554:VAL:HG13	1:B:555:GLY:H	1.83	0.44
1:C:339:GLU:O	1:C:340:THR:C	2.60	0.44
1:D:38:SER:OG	1:D:40:PRO:HD3	2.18	0.44
1:D:427:THR:OG1	1:D:578:THR:HG22	2.17	0.44
1:B:470:PHE:CG	1:B:525:LEU:HD22	2.53	0.43
1:A:478:PHE:CE1	1:A:498:ILE:HA	2.54	0.43
1:B:260:GLU:HB2	1:B:262:TYR:CE2	2.53	0.43
1:C:43:ASN:HB2	1:C:69:CYS:O	2.18	0.43
1:C:133:HIS:CD2	1:C:147:TYR:HE2	2.36	0.43
1:C:463:LEU:O	1:C:467:ARG:HG3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:VAL:H	1:B:459:LYS:HB3	1.82	0.43
1:A:479:GLU:CD	1:A:479:GLU:N	2.75	0.43
1:B:126:SER:HA	1:B:127:PRO:C	2.43	0.43
1:B:464:ASN:ND2	1:B:475:TYR:H	2.15	0.43
1:B:565:GLN:O	1:B:569:CYS:HB2	2.19	0.43
1:C:133:HIS:HD2	1:C:147:TYR:CZ	2.37	0.43
1:C:155:VAL:H	1:C:459:LYS:HB3	1.83	0.43
1:C:287:VAL:HG23	1:C:288:GLY:N	2.34	0.43
1:D:195:ASN:HB3	1:D:582:VAL:HB	1.99	0.43
1:D:388:HIS:N	1:D:389:PRO:CD	2.81	0.43
1:A:423:VAL:CG1	1:A:578:THR:HG23	2.48	0.43
1:B:485:LYS:HD2	1:B:489:ALA:HB2	2.00	0.43
1:C:124:ILE:N	1:C:124:ILE:HD12	2.34	0.43
1:C:434:VAL:H	1:C:517:ILE:HG13	1.82	0.43
1:D:113:MET:HE2	1:D:360:LYS:O	2.19	0.43
1:D:396:ASN:HD22	1:D:401:GLU:HB3	1.84	0.43
1:A:294:LEU:HD23	1:A:409:TYR:HB3	2.00	0.43
1:B:122:TYR:CD1	1:B:122:TYR:C	2.97	0.43
1:B:245:ARG:HH22	1:B:326:GLU:HG2	1.84	0.43
1:B:336:LEU:N	1:B:336:LEU:CD1	2.81	0.43
1:B:453:ASP:O	1:B:456:ARG:HB2	2.19	0.43
1:C:43:ASN:O	1:C:44:ARG:CB	2.67	0.43
1:C:206:THR:HG21	1:C:385:TYR:CE1	2.54	0.43
1:D:178:LEU:HA	1:D:182:LEU:HD12	2.01	0.43
1:D:500:VAL:O	1:D:500:VAL:HG12	2.18	0.43
1:B:129:THR:CG2	1:B:137:LYS:HD2	2.49	0.43
1:B:178:LEU:HA	1:B:182:LEU:HD12	2.00	0.43
1:B:213:ASP:OD1	1:B:215:LYS:HB2	2.18	0.43
1:C:316:LEU:HD12	1:C:316:LEU:HA	1.93	0.43
1:C:428:ARG:HA	1:C:583:GLN:OE1	2.19	0.43
1:C:478:PHE:HE2	1:C:495:TYR:HB2	1.81	0.43
1:D:410:ASN:OD1	1:D:413:ILE:HG13	2.19	0.43
1:D:487:MET:HA	1:D:490:GLU:HB3	2.00	0.43
1:D:495:TYR:CE2	1:D:501:MET:SD	3.11	0.43
1:B:178:LEU:O	1:B:182:LEU:HB2	2.18	0.43
1:B:352:LEU:HD21	1:B:387:TRP:CH2	2.53	0.43
1:C:145:LEU:HD23	1:C:376:ARG:HE	1.84	0.43
1:D:477:SER:HB2	1:D:479:GLU:CG	2.35	0.43
1:A:127:PRO:O	1:B:543:GLN:NE2	2.52	0.43
1:B:403:SER:O	1:B:404:PHE:C	2.61	0.43
1:C:113:MET:HE3	1:C:117:LEU:CD1	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:43:ASN:HB2	1:D:69:CYS:O	2.19	0.43
1:A:120:ARG:NH1	4:A:701:IMN:O2	2.42	0.42
1:A:353:SER:HA	4:A:701:IMN:C3	2.49	0.42
1:A:388:HIS:HB3	1:A:444:VAL:HG21	2.01	0.42
1:A:396:ASN:ND2	1:A:401:GLU:HB3	2.34	0.42
1:B:522:MET:HE2	1:B:522:MET:HB2	1.96	0.42
1:C:89:VAL:HG12	1:C:93:LEU:HD23	2.01	0.42
1:C:273:MET:HG3	1:C:285:PHE:O	2.19	0.42
1:A:500:VAL:O	1:A:500:VAL:HG12	2.18	0.42
1:D:192:GLN:HG3	1:D:516:ALA:HA	2.02	0.42
1:B:273:MET:SD	1:B:290:GLU:HA	2.60	0.42
1:B:513:ARG:HG3	1:B:518:PHE:O	2.20	0.42
1:C:260:GLU:HB2	1:C:262:TYR:CE2	2.54	0.42
1:A:149:THR:O	1:A:378:ALA:HA	2.19	0.42
1:A:252:LEU:O	1:A:310:GLN:NE2	2.52	0.42
1:A:335:ILE:HA	1:A:559:ILE:HD11	2.01	0.42
1:B:178:LEU:HD13	1:B:183:LEU:HD11	2.02	0.42
1:B:396:ASN:ND2	1:B:401:GLU:HB3	2.34	0.42
1:C:137:LYS:O	1:C:138:SER:O	2.36	0.42
1:C:553:GLU:O	1:C:557:LYS:NZ	2.47	0.42
1:D:245:ARG:NH2	1:D:247:PHE:CE1	2.87	0.42
1:D:333:LYS:O	1:D:337:ILE:HG13	2.19	0.42
1:A:107:PHE:CD1	1:A:107:PHE:N	2.85	0.42
1:A:126:SER:HB2	1:A:532:LYS:NZ	2.35	0.42
1:A:145:LEU:HD23	1:A:376:ARG:HH21	1.85	0.42
1:A:229:ASP:OD1	1:A:231:ASN:HB3	2.18	0.42
1:A:372:GLN:OE1	1:A:374:GLN:NE2	2.53	0.42
1:B:475:TYR:HA	1:B:480:GLU:OE1	2.19	0.42
1:B:568:ILE:HG23	1:B:572:VAL:CG2	2.50	0.42
3:B:682:HEM:HBC2	3:B:682:HEM:CHD	2.41	0.42
1:C:74:PHE:HA	1:C:77:ARG:HD2	2.01	0.42
1:D:339:GLU:HG2	1:D:562:ALA:HB2	2.02	0.42
1:A:134:TYR:CD2	1:A:136:TYR:CE1	3.05	0.42
1:A:149:THR:HG21	1:A:376:ARG:HB3	2.01	0.42
1:A:254:TYR:HA	1:A:264:PRO:HD3	2.02	0.42
1:B:124:ILE:HD11	1:B:528:PRO:CB	2.49	0.42
1:B:131:ASN:ND2	1:B:147:TYR:CD2	2.87	0.42
1:C:423:VAL:HG13	1:C:578:THR:HG23	2.00	0.42
1:D:206:THR:HB	1:D:210:PHE:CD2	2.54	0.42
1:D:564:ILE:HG12	1:D:580:PHE:CE1	2.54	0.42
1:A:43:ASN:HB2	1:A:69:CYS:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:THR:O	1:A:91:TYR:HB3	2.19	0.42
1:A:479:GLU:HA	1:A:488:ALA:CB	2.50	0.42
1:B:287:VAL:HG11	1:B:299:MET:SD	2.60	0.42
1:B:352:LEU:HD12	1:B:352:LEU:HA	1.73	0.42
1:B:447:VAL:O	1:B:450:ALA:HB3	2.19	0.42
1:B:487:MET:HA	1:B:490:GLU:HB3	2.02	0.42
1:C:134:TYR:CD2	1:C:136:TYR:CE1	3.06	0.42
1:C:495:TYR:CD2	1:C:501:MET:SD	3.13	0.42
1:D:287:VAL:HG23	1:D:288:GLY:N	2.35	0.42
1:A:43:ASN:O	1:A:44:ARG:CB	2.68	0.42
1:A:230:LEU:HG	1:A:337:ILE:CG1	2.49	0.42
1:C:403:SER:OG	1:C:406:GLN:HG3	2.19	0.42
1:C:526:GLY:HA3	4:C:701:IMN:C14	2.50	0.42
1:D:96:PHE:CD1	1:D:96:PHE:N	2.88	0.42
1:D:234:TYR:OH	1:D:336:LEU:HD22	2.20	0.42
1:A:206:THR:HG21	1:A:385:TYR:CE1	2.55	0.42
1:B:465:GLU:O	1:B:468:LYS:HB2	2.19	0.42
1:B:479:GLU:HA	1:B:488:ALA:CB	2.50	0.42
1:C:127:PRO:HD2	1:D:543:GLN:HE22	1.84	0.42
1:B:34:ASN:HB3	1:B:37:CYS:SG	2.60	0.42
1:B:149:THR:O	1:B:378:ALA:HA	2.19	0.42
1:B:391:LEU:HD21	3:B:682:HEM:HAB	2.02	0.42
1:C:65:TYR:HD1	1:C:73:GLU:OE2	2.02	0.42
1:D:281:GLU:C	1:D:283:LEU:N	2.76	0.42
1:D:465:GLU:O	1:D:468:LYS:HB2	2.19	0.42
1:A:495:TYR:O	1:A:497:ASP:N	2.53	0.41
1:C:557:LYS:O	1:C:558:ILE:C	2.62	0.41
1:D:124:ILE:HD12	1:D:124:ILE:H	1.85	0.41
1:A:253:LYS:HD3	1:A:269:THR:HG22	2.01	0.41
1:A:459:LYS:HA	1:A:459:LYS:HE2	2.02	0.41
1:B:341:ILE:O	1:B:345:ILE:HG13	2.20	0.41
1:B:384:LEU:HD23	1:B:507:LEU:HD11	2.02	0.41
1:C:574:GLY:C	1:C:576:PRO:HD3	2.42	0.41
1:A:124:ILE:HG22	1:A:125:ASP:N	2.35	0.41
1:A:298:LEU:HD12	1:A:298:LEU:HA	1.92	0.41
1:A:322:GLU:HG2	1:B:51:GLY:C	2.45	0.41
1:B:335:ILE:HA	1:B:559:ILE:HD11	2.02	0.41
1:C:150:ARG:NH1	1:C:150:ARG:CG	2.83	0.41
1:C:513:ARG:O	1:C:514:PRO:C	2.63	0.41
1:D:106:PRO:O	1:D:107:PHE:C	2.63	0.41
1:D:339:GLU:O	1:D:340:THR:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:391:LEU:HD21	3:A:682:HEM:HAB	2.02	0.41
1:A:485:LYS:HD3	1:A:488:ALA:HB3	2.02	0.41
1:B:113:MET:O	1:B:116:VAL:HG22	2.20	0.41
1:C:161:THR:HG21	1:C:165:VAL:O	2.20	0.41
1:C:522:MET:HE2	1:C:522:MET:HB2	1.87	0.41
1:D:206:THR:HG21	1:D:385:TYR:CE1	2.55	0.41
1:D:269:THR:O	1:D:270:GLN:CB	2.67	0.41
1:D:568:ILE:O	1:D:572:VAL:HG23	2.19	0.41
1:A:341:ILE:O	1:A:345:ILE:HG13	2.21	0.41
1:B:423:VAL:CG1	1:B:578:THR:HG23	2.50	0.41
1:C:513:ARG:HG3	1:C:518:PHE:O	2.19	0.41
1:D:477:SER:O	1:D:480:GLU:N	2.53	0.41
1:A:46:GLU:OE2	1:A:137:LYS:HE3	2.21	0.41
1:A:106:PRO:O	1:A:109:ARG:N	2.54	0.41
1:A:129:THR:CG2	1:A:137:LYS:HD2	2.51	0.41
1:A:169:LYS:O	1:A:170:GLU:HG3	2.21	0.41
1:A:213:ASP:OD1	1:A:215:LYS:HB2	2.21	0.41
1:B:525:LEU:N	1:B:525:LEU:CD2	2.83	0.41
1:C:106:PRO:O	1:C:109:ARG:N	2.54	0.41
1:C:487:MET:HA	1:C:490:GLU:HB3	2.01	0.41
1:D:254:TYR:CD1	1:D:254:TYR:C	2.99	0.41
1:A:133:HIS:HD2	1:A:147:TYR:OH	2.03	0.41
1:A:137:LYS:O	1:A:138:SER:O	2.39	0.41
1:A:495:TYR:CE2	1:A:501:MET:SD	3.13	0.41
1:B:184:ARG:NH1	1:B:441:PRO:HG3	2.36	0.41
1:C:131:ASN:OD1	1:C:131:ASN:C	2.63	0.41
1:C:209:PHE:HB2	1:C:377:ILE:HG21	2.02	0.41
1:D:398:GLU:HG2	1:D:421:GLN:NE2	2.36	0.41
1:D:512:PRO:HB3	1:D:516:ALA:O	2.21	0.41
1:A:209:PHE:HB2	1:A:377:ILE:HG21	2.03	0.41
1:A:248:LYS:HE2	1:A:248:LYS:HB2	1.95	0.41
1:A:339:GLU:O	1:A:342:LYS:HB3	2.20	0.41
1:A:366:LEU:HA	1:A:369:GLN:HG2	2.02	0.41
1:A:557:LYS:N	1:A:557:LYS:HE3	2.35	0.41
1:B:105(A):ILE:O	1:B:106:PRO:C	2.64	0.41
1:B:181:VAL:HB	1:B:509:VAL:HG21	2.02	0.41
1:B:339:GLU:O	1:B:340:THR:C	2.63	0.41
1:C:76:THR:O	1:C:80:LEU:HG	2.20	0.41
1:D:105(A):ILE:O	1:D:106:PRO:C	2.64	0.41
1:D:283:LEU:CD2	1:D:415:LEU:HD12	2.50	0.41
1:D:352:LEU:HD12	1:D:352:LEU:HA	1.76	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:423:VAL:CG1	1:D:578:THR:HG23	2.50	0.41
1:D:522:MET:HE2	1:D:522:MET:HB2	1.93	0.41
1:A:50:THR:HG21	1:A:56:LYS:HB2	2.02	0.41
1:A:52:PHE:HA	1:B:322:GLU:HG2	2.02	0.41
1:A:74:PHE:HD1	1:A:77:ARG:HH11	1.66	0.41
1:A:96:PHE:O	1:A:97:LYS:C	2.64	0.41
1:A:464:ASN:ND2	1:A:475:TYR:H	2.16	0.41
1:A:522:MET:HE2	1:A:522:MET:HB2	1.94	0.41
1:B:145:LEU:HD23	1:B:376:ARG:HE	1.86	0.41
1:C:180:LYS:HD3	1:C:490:GLU:OE1	2.21	0.41
1:C:387:TRP:HB2	3:C:682:HEM:CAC	2.49	0.41
1:C:445:GLN:HG3	1:C:446:ALA:N	2.35	0.41
1:C:469:ARG:HG3	1:C:470:PHE:CE1	2.56	0.41
1:C:500:VAL:HG12	1:C:500:VAL:O	2.21	0.41
1:C:554:VAL:HG13	1:C:555:GLY:H	1.86	0.41
1:D:91:TYR:O	1:D:95:HIS:CD2	2.74	0.41
1:D:341:ILE:O	1:D:345:ILE:HG13	2.21	0.41
1:D:368:ASN:OD1	1:D:369:GLN:NE2	2.50	0.41
1:D:389:PRO:O	1:D:432:GLY:HA3	2.21	0.41
1:D:391:LEU:HB2	1:D:441:PRO:HG2	2.02	0.41
1:D:557:LYS:HD3	1:D:557:LYS:HA	1.82	0.41
1:D:577:PHE:CD1	1:D:577:PHE:C	2.99	0.41
1:A:96:PHE:CD1	1:A:96:PHE:N	2.88	0.41
1:A:391:LEU:HB2	1:A:441:PRO:HG2	2.03	0.41
1:A:552:GLY:HA2	1:B:48:MET:HE1	2.03	0.41
1:C:194:SER:OG	1:C:351:HIS:HE1	2.04	0.41
1:A:243:LYS:HB3	1:A:271:VAL:HG12	2.03	0.40
1:B:126:SER:HB2	1:B:532:LYS:NZ	2.35	0.40
1:B:134:TYR:CD2	1:B:136:TYR:CE1	3.06	0.40
1:B:152:LEU:HD23	1:B:152:LEU:HA	1.93	0.40
1:B:423:VAL:HG11	1:B:576:PRO:HB2	2.03	0.40
1:B:464:ASN:HD21	1:B:475:TYR:N	2.19	0.40
1:C:152:LEU:HD23	1:C:152:LEU:HA	1.95	0.40
1:C:352:LEU:HD12	1:C:352:LEU:HA	1.87	0.40
1:C:485:LYS:HD3	1:C:488:ALA:HB3	2.02	0.40
1:D:38:SER:O	1:D:39:ASN:C	2.63	0.40
1:D:254:TYR:HD2	1:D:310:GLN:HE21	1.68	0.40
1:D:316:LEU:HD12	1:D:316:LEU:HA	1.82	0.40
1:D:479:GLU:HB3	1:D:485:LYS:CE	2.51	0.40
1:A:339:GLU:O	1:A:340:THR:C	2.65	0.40
1:B:500:VAL:O	1:B:500:VAL:HG12	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:198:PHE:HD1	1:D:199:ALA:N	2.19	0.40
1:D:579:SER:OG	1:D:580:PHE:N	2.54	0.40
1:A:93:LEU:HD13	1:A:93:LEU:HA	1.93	0.40
1:A:150:ARG:NH2	1:A:154:PRO:HB3	2.35	0.40
1:A:276:PRO:HA	1:A:277:PRO:HD2	1.86	0.40
1:A:287:VAL:HG11	1:A:299:MET:SD	2.61	0.40
1:A:380:GLU:H	1:A:380:GLU:CD	2.30	0.40
1:B:149:THR:HG21	1:B:376:ARG:HB3	2.02	0.40
1:C:124:ILE:HD11	1:C:528:PRO:CB	2.51	0.40
1:C:464:ASN:ND2	1:C:475:TYR:H	2.19	0.40
1:C:518:PHE:HB3	1:C:522:MET:HB3	2.03	0.40
1:D:42:GLN:HE22	1:D:468:LYS:HE2	1.85	0.40
1:D:352:LEU:HD21	1:D:387:TRP:HH2	1.84	0.40
1:A:92:ILE:HG13	1:A:92:ILE:H	1.70	0.40
1:A:260:GLU:HB2	1:A:262:TYR:CE2	2.54	0.40
1:B:43:ASN:HB2	1:B:69:CYS:O	2.21	0.40
1:B:150:ARG:HG3	1:B:380:GLU:OE2	2.22	0.40
1:B:213:ASP:HB2	1:B:222:ARG:HG3	2.04	0.40
1:B:276:PRO:HD2	1:B:279:ILE:CD1	2.52	0.40
1:C:121:SER:O	1:C:123:LEU:N	2.55	0.40
1:C:345:ILE:HG12	1:C:534:LEU:HD23	2.02	0.40
1:C:568:ILE:HG23	1:C:572:VAL:CG2	2.50	0.40
1:D:500:VAL:O	1:D:500:VAL:CG1	2.70	0.40
1:B:234:TYR:HE1	1:B:309:HIS:CE1	2.39	0.40
1:B:276:PRO:HG2	1:B:409:TYR:CG	2.57	0.40
1:B:294:LEU:HD23	1:B:409:TYR:HB3	2.03	0.40
1:B:464:ASN:OD1	1:B:499:ASP:HA	2.22	0.40
1:C:39:ASN:N	1:C:40:PRO:CD	2.84	0.40
1:C:512:PRO:HB3	1:C:516:ALA:O	2.21	0.40
1:D:415:LEU:HA	1:D:415:LEU:HD23	1.84	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	550/587 (94%)	456 (83%)	74 (14%)	20 (4%)	2	11
1	B	550/587 (94%)	455 (83%)	74 (14%)	21 (4%)	2	10
1	C	550/587 (94%)	454 (82%)	78 (14%)	18 (3%)	3	13
1	D	550/587 (94%)	451 (82%)	76 (14%)	23 (4%)	2	9
All	All	2200/2348 (94%)	1816 (82%)	302 (14%)	82 (4%)	2	11

All (82) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	138	SER
1	A	282	ASN
1	A	398	GLU
1	A	514	PRO
1	B	138	SER
1	B	282	ASN
1	B	398	GLU
1	B	514	PRO
1	C	138	SER
1	C	282	ASN
1	C	398	GLU
1	C	422	PHE
1	C	514	PRO
1	D	138	SER
1	D	282	ASN
1	D	398	GLU
1	D	422	PHE
1	D	514	PRO
1	D	575	CYS
1	A	82	LEU
1	A	419	LEU
1	A	422	PHE
1	A	575	CYS
1	B	419	LEU
1	B	422	PHE
1	B	575	CYS
1	C	43	ASN
1	D	419	LEU
1	A	169	LYS
1	B	43	ASN

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Mol	Chain	Res	Type
1	B	82	LEU
1	B	438	ARG
1	C	82	LEU
1	C	130	TYR
1	C	280	PRO
1	C	419	LEU
1	D	82	LEU
1	D	438	ARG
1	A	43	ASN
1	A	280	PRO
1	A	325	ASP
1	B	280	PRO
1	B	325	ASP
1	B	348	TYR
1	C	325	ASP
1	C	438	ARG
1	C	575	CYS
1	D	43	ASN
1	D	280	PRO
1	D	325	ASP
1	D	348	TYR
1	A	106	PRO
1	A	130	TYR
1	A	348	TYR
1	A	418	GLY
1	A	438	ARG
1	B	169	LYS
1	B	496	SER
1	C	416	GLU
1	C	496	SER
1	D	106	PRO
1	D	130	TYR
1	D	418	GLY
1	D	429	GLN
1	D	496	SER
1	B	106	PRO
1	B	429	GLN
1	C	106	PRO
1	A	444	VAL
1	B	287	VAL
1	B	444	VAL
1	C	444	VAL

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Mol	Chain	Res	Type
1	D	444	VAL
1	C	418	GLY
1	D	500	VAL
1	A	132	VAL
1	B	418	GLY
1	D	132	VAL
1	D	528	PRO
1	A	500	VAL
1	B	430	ILE
1	D	287	VAL

5.3.2 Protein sidechains [i](#)

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	493/525 (94%)	423 (86%)	70 (14%)	3	11
1	B	493/525 (94%)	421 (85%)	72 (15%)	3	10
1	C	493/525 (94%)	423 (86%)	70 (14%)	3	11
1	D	493/525 (94%)	426 (86%)	67 (14%)	3	12
All	All	1972/2100 (94%)	1693 (86%)	279 (14%)	3	11

All (279) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	ARG
1	A	48	MET
1	A	49	SER
1	A	70	THR
1	A	71	THR
1	A	73	GLU
1	A	75	LEU
1	A	82	LEU
1	A	83	LYS
1	A	93	LEU
1	A	94	THR

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Mol	Chain	Res	Type
1	A	103	VAL
1	A	111	LEU
1	A	116	VAL
1	A	117	LEU
1	A	120	ARG
1	A	122	TYR
1	A	126	SER
1	A	137	LYS
1	A	150	ARG
1	A	162	PRO
1	A	170	GLU
1	A	173	ASP
1	A	178	LEU
1	A	186	GLU
1	A	189	PRO
1	A	209	PHE
1	A	230	LEU
1	A	231	ASN
1	A	238	LEU
1	A	241	GLN
1	A	248	LYS
1	A	252	LEU
1	A	270	GLN
1	A	271	VAL
1	A	282	ASN
1	A	289	GLN
1	A	290	GLU
1	A	291	VAL
1	A	298	LEU
1	A	300	MET
1	A	310	GLN
1	A	311	ARG
1	A	316	LEU
1	A	322	GLU
1	A	377	ILE
1	A	385	TYR
1	A	389	PRO
1	A	398	GLU
1	A	400	GLN
1	A	405	LYS
1	A	412	SER
1	A	420	THR

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Mol	Chain	Res	Type
1	A	428	ARG
1	A	441	PRO
1	A	469	ARG
1	A	476	THR
1	A	479	GLU
1	A	484	GLU
1	A	485	LYS
1	A	486	GLU
1	A	514	PRO
1	A	525	LEU
1	A	532	LYS
1	A	534	LEU
1	A	539	ILE
1	A	543	GLN
1	A	557	LYS
1	A	578	THR
1	A	583	GLN
1	B	39	ASN
1	B	44	ARG
1	B	49	SER
1	B	70	THR
1	B	71	THR
1	B	73	GLU
1	B	75	LEU
1	B	82	LEU
1	B	83	LYS
1	B	93	LEU
1	B	94	THR
1	B	103	VAL
1	B	106	PRO
1	B	111	LEU
1	B	116	VAL
1	B	117	LEU
1	B	120	ARG
1	B	122	TYR
1	B	126	SER
1	B	137	LYS
1	B	138	SER
1	B	150	ARG
1	B	162	PRO
1	B	170	GLU
1	B	178	LEU

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Mol	Chain	Res	Type
1	B	186	GLU
1	B	189	PRO
1	B	209	PHE
1	B	230	LEU
1	B	231	ASN
1	B	236	GLU
1	B	238	LEU
1	B	241	GLN
1	B	248	LYS
1	B	252	LEU
1	B	271	VAL
1	B	280	PRO
1	B	281	GLU
1	B	282	ASN
1	B	289	GLN
1	B	290	GLU
1	B	291	VAL
1	B	298	LEU
1	B	300	MET
1	B	310	GLN
1	B	311	ARG
1	B	316	LEU
1	B	322	GLU
1	B	377	ILE
1	B	385	TYR
1	B	389	PRO
1	B	398	GLU
1	B	400	GLN
1	B	412	SER
1	B	420	THR
1	B	428	ARG
1	B	441	PRO
1	B	469	ARG
1	B	476	THR
1	B	479	GLU
1	B	484	GLU
1	B	485	LYS
1	B	486	GLU
1	B	514	PRO
1	B	525	LEU
1	B	532	LYS
1	B	534	LEU

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Mol	Chain	Res	Type
1	B	539	ILE
1	B	543	GLN
1	B	557	LYS
1	B	578	THR
1	B	583	GLN
1	C	44	ARG
1	C	48	MET
1	C	49	SER
1	C	70	THR
1	C	71	THR
1	C	73	GLU
1	C	75	LEU
1	C	82	LEU
1	C	83	LYS
1	C	93	LEU
1	C	94	THR
1	C	111	LEU
1	C	116	VAL
1	C	117	LEU
1	C	120	ARG
1	C	122	TYR
1	C	126	SER
1	C	137	LYS
1	C	150	ARG
1	C	162	PRO
1	C	170	GLU
1	C	173	ASP
1	C	178	LEU
1	C	186	GLU
1	C	189	PRO
1	C	209	PHE
1	C	230	LEU
1	C	231	ASN
1	C	236	GLU
1	C	238	LEU
1	C	241	GLN
1	C	248	LYS
1	C	252	LEU
1	C	271	VAL
1	C	281	GLU
1	C	282	ASN
1	C	289	GLN

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Mol	Chain	Res	Type
1	C	290	GLU
1	C	291	VAL
1	C	298	LEU
1	C	300	MET
1	C	310	GLN
1	C	311	ARG
1	C	316	LEU
1	C	322	GLU
1	C	352	LEU
1	C	377	ILE
1	C	385	TYR
1	C	389	PRO
1	C	398	GLU
1	C	400	GLN
1	C	405	LYS
1	C	412	SER
1	C	420	THR
1	C	441	PRO
1	C	469	ARG
1	C	476	THR
1	C	479	GLU
1	C	484	GLU
1	C	485	LYS
1	C	486	GLU
1	C	514	PRO
1	C	525	LEU
1	C	532	LYS
1	C	534	LEU
1	C	539	ILE
1	C	543	GLN
1	C	557	LYS
1	C	578	THR
1	C	583	GLN
1	D	44	ARG
1	D	49	SER
1	D	70	THR
1	D	71	THR
1	D	73	GLU
1	D	75	LEU
1	D	82	LEU
1	D	83	LYS
1	D	93	LEU

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Mol	Chain	Res	Type
1	D	94	THR
1	D	106	PRO
1	D	111	LEU
1	D	116	VAL
1	D	117	LEU
1	D	120	ARG
1	D	122	TYR
1	D	126	SER
1	D	137	LYS
1	D	150	ARG
1	D	162	PRO
1	D	170	GLU
1	D	173	ASP
1	D	174	SER
1	D	178	LEU
1	D	186	GLU
1	D	209	PHE
1	D	230	LEU
1	D	231	ASN
1	D	236	GLU
1	D	238	LEU
1	D	241	GLN
1	D	248	LYS
1	D	252	LEU
1	D	282	ASN
1	D	289	GLN
1	D	290	GLU
1	D	291	VAL
1	D	298	LEU
1	D	300	MET
1	D	310	GLN
1	D	311	ARG
1	D	316	LEU
1	D	322	GLU
1	D	377	ILE
1	D	385	TYR
1	D	389	PRO
1	D	398	GLU
1	D	400	GLN
1	D	405	LYS
1	D	412	SER
1	D	420	THR

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Mol	Chain	Res	Type
1	D	428	ARG
1	D	441	PRO
1	D	469	ARG
1	D	476	THR
1	D	479	GLU
1	D	484	GLU
1	D	485	LYS
1	D	486	GLU
1	D	514	PRO
1	D	525	LEU
1	D	532	LYS
1	D	534	LEU
1	D	543	GLN
1	D	557	LYS
1	D	578	THR
1	D	583	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (64) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	A	43	ASN
1	A	95	HIS
1	A	105	ASN
1	A	133	HIS
1	A	232	HIS
1	A	320	HIS
1	A	369	GLN
1	A	372	GLN
1	A	374	GLN
1	A	382	ASN
1	A	396	ASN
1	A	411	ASN
1	A	454	GLN
1	B	42	GLN
1	B	54	GLN
1	B	95	HIS
1	B	133	HIS
1	B	232	HIS
1	B	278	HIS
1	B	282	ASN
1	B	320	HIS

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Mol	Chain	Res	Type
1	B	356	HIS
1	B	369	GLN
1	B	372	GLN
1	B	374	GLN
1	B	382	ASN
1	B	396	ASN
1	B	411	ASN
1	B	454	GLN
1	B	464	ASN
1	C	43	ASN
1	C	105	ASN
1	C	133	HIS
1	C	214	HIS
1	C	232	HIS
1	C	242	HIS
1	C	278	HIS
1	C	282	ASN
1	C	320	HIS
1	C	351	HIS
1	C	369	GLN
1	C	372	GLN
1	C	374	GLN
1	C	382	ASN
1	C	396	ASN
1	C	411	ASN
1	C	454	GLN
1	C	464	ASN
1	C	543	GLN
1	D	42	GLN
1	D	43	ASN
1	D	133	HIS
1	D	232	HIS
1	D	278	HIS
1	D	282	ASN
1	D	320	HIS
1	D	356	HIS
1	D	372	GLN
1	D	374	GLN
1	D	382	ASN
1	D	396	ASN
1	D	411	ASN
1	D	454	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

20 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	NAG	D	671	1	14,14,15	0.65	1 (7%)	17,19,21	1.17	3 (17%)
3	HEM	A	682	1	50,50,50	1.30	5 (10%)	67,82,82	0.91	1 (1%)
4	IMN	D	701	-	27,27,27	2.30	10 (37%)	39,39,39	1.29	5 (12%)
2	NAG	D	661	1	14,14,15	0.76	0	17,19,21	0.88	1 (5%)
2	NAG	C	661	1	14,14,15	0.81	0	17,19,21	0.84	2 (11%)
2	NAG	D	681	1	14,14,15	0.80	0	17,19,21	0.96	0
2	NAG	B	671	1	14,14,15	0.70	0	17,19,21	1.09	2 (11%)
2	NAG	A	661	1	14,14,15	0.59	0	17,19,21	0.81	1 (5%)
4	IMN	A	701	-	27,27,27	2.09	10 (37%)	39,39,39	1.31	4 (10%)
3	HEM	B	682	1	50,50,50	1.23	4 (8%)	67,82,82	1.03	3 (4%)
3	HEM	D	682	1	50,50,50	1.26	5 (10%)	67,82,82	0.94	1 (1%)
4	IMN	B	701	-	27,27,27	2.15	12 (44%)	39,39,39	1.25	5 (12%)
2	NAG	A	681	1	14,14,15	0.44	0	17,19,21	0.88	1 (5%)
2	NAG	A	671	1	14,14,15	0.59	0	17,19,21	1.02	0
4	IMN	C	701	-	27,27,27	2.09	11 (40%)	39,39,39	1.32	5 (12%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	C	681	1	14,14,15	0.49	0	17,19,21	0.80	0
2	NAG	C	671	1	14,14,15	0.85	0	17,19,21	1.23	2 (11%)
3	HEM	C	682	1	50,50,50	1.34	6 (12%)	67,82,82	0.90	1 (1%)
2	NAG	B	661	1	14,14,15	0.82	0	17,19,21	0.86	1 (5%)
2	NAG	B	681	1	14,14,15	0.47	0	17,19,21	0.87	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	D	671	1	-	2/6/23/26	0/1/1/1
3	HEM	A	682	1	-	4/14/54/54	-
4	IMN	D	701	-	-	9/14/14/14	0/3/3/3
2	NAG	D	661	1	-	2/6/23/26	0/1/1/1
2	NAG	C	661	1	-	2/6/23/26	0/1/1/1
2	NAG	D	681	1	-	0/6/23/26	0/1/1/1
2	NAG	B	671	1	-	2/6/23/26	0/1/1/1
2	NAG	A	661	1	-	2/6/23/26	0/1/1/1
4	IMN	A	701	-	-	10/14/14/14	0/3/3/3
3	HEM	B	682	1	-	3/14/54/54	-
3	HEM	D	682	1	-	2/14/54/54	-
4	IMN	B	701	-	-	9/14/14/14	0/3/3/3
2	NAG	A	681	1	-	2/6/23/26	0/1/1/1
2	NAG	A	671	1	-	2/6/23/26	0/1/1/1
4	IMN	C	701	-	-	9/14/14/14	0/3/3/3
2	NAG	C	681	1	-	0/6/23/26	0/1/1/1
2	NAG	C	671	1	-	2/6/23/26	0/1/1/1
3	HEM	C	682	1	-	4/14/54/54	-
2	NAG	B	661	1	-	2/6/23/26	0/1/1/1
2	NAG	B	681	1	-	0/6/23/26	0/1/1/1

All (64) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	701	IMN	C5-C	5.17	1.47	1.39
4	A	701	IMN	C1-C7	4.47	1.52	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	682	HEM	CBB-CAB	4.38	1.51	1.30
4	D	701	IMN	C2-C3	4.22	1.45	1.39
4	D	701	IMN	C5-C4	4.13	1.45	1.38
4	B	701	IMN	C5-C	3.90	1.45	1.39
4	D	701	IMN	C15-C14	3.76	1.44	1.38
3	A	682	HEM	CBB-CAB	3.76	1.48	1.30
3	D	682	HEM	CAC-C3C	-3.73	1.37	1.47
3	D	682	HEM	CBB-CAB	3.71	1.48	1.30
4	D	701	IMN	C9-N	-3.70	1.32	1.41
4	C	701	IMN	C1-C7	3.69	1.50	1.44
4	C	701	IMN	C4-C3	3.64	1.45	1.38
4	A	701	IMN	C9-N	-3.64	1.33	1.41
4	B	701	IMN	C4-C3	3.51	1.45	1.38
3	B	682	HEM	CBB-CAB	3.50	1.47	1.30
3	A	682	HEM	CBC-CAC	3.43	1.46	1.30
3	A	682	HEM	CAC-C3C	-3.41	1.38	1.47
4	A	701	IMN	C5-C	3.35	1.44	1.39
3	C	682	HEM	CBC-CAC	3.34	1.46	1.30
4	B	701	IMN	C1-C7	3.33	1.50	1.44
4	C	701	IMN	C5-C	3.31	1.44	1.39
4	B	701	IMN	C15-C14	3.31	1.44	1.38
4	D	701	IMN	C1-C7	3.27	1.50	1.44
3	A	682	HEM	CAB-C3B	-3.25	1.38	1.47
4	C	701	IMN	C15-C14	3.24	1.44	1.38
3	D	682	HEM	CBC-CAC	3.20	1.45	1.30
4	B	701	IMN	C9-N	-3.18	1.34	1.41
4	A	701	IMN	C4-C3	3.18	1.44	1.38
4	C	701	IMN	C9-N	-3.17	1.34	1.41
4	B	701	IMN	C8-N	-3.14	1.34	1.40
3	C	682	HEM	CAB-C3B	-3.14	1.39	1.47
3	B	682	HEM	CAC-C3C	-3.12	1.39	1.47
4	C	701	IMN	C5-C4	3.10	1.43	1.38
4	A	701	IMN	C15-C14	3.03	1.43	1.38
3	D	682	HEM	CAB-C3B	-3.02	1.39	1.47
3	B	682	HEM	CAB-C3B	-2.99	1.39	1.47
4	A	701	IMN	C5-C4	2.97	1.43	1.38
3	C	682	HEM	CAC-C3C	-2.88	1.39	1.47
4	D	701	IMN	C4-C3	2.87	1.44	1.38
4	B	701	IMN	C2-C3	2.85	1.43	1.39
3	B	682	HEM	CBC-CAC	2.82	1.44	1.30
4	C	701	IMN	C2-C3	2.79	1.43	1.39
4	C	701	IMN	C17-C7	2.73	1.56	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	682	HEM	CBA-CGA	2.69	1.56	1.50
4	B	701	IMN	C5-C4	2.61	1.43	1.38
4	A	701	IMN	C17-C7	2.59	1.56	1.50
4	A	701	IMN	O-C3	2.51	1.42	1.37
3	C	682	HEM	C1B-NB	-2.49	1.35	1.40
4	A	701	IMN	C2-C3	2.41	1.42	1.39
4	B	701	IMN	C2-C1	2.37	1.43	1.39
4	D	701	IMN	C8-N	-2.35	1.35	1.40
4	B	701	IMN	C12-C11	2.26	1.42	1.38
4	C	701	IMN	C10-C9	2.22	1.53	1.50
4	D	701	IMN	C2-C1	2.19	1.43	1.39
4	A	701	IMN	C2-C1	2.15	1.43	1.39
4	B	701	IMN	O-C3	2.13	1.41	1.37
3	D	682	HEM	CBA-CGA	2.10	1.55	1.50
4	B	701	IMN	C10-C9	2.09	1.53	1.50
4	C	701	IMN	C12-C11	2.08	1.42	1.38
2	D	671	NAG	C1-C2	2.07	1.55	1.52
4	D	701	IMN	C1-C	2.05	1.43	1.41
4	C	701	IMN	C8-N	-2.04	1.36	1.40
3	C	682	HEM	O2D-CGD	-2.02	1.24	1.30

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	682	HEM	C3B-C4B-NB	3.73	112.15	109.47
3	C	682	HEM	C3B-C4B-NB	3.44	111.94	109.47
3	D	682	HEM	C3B-C4B-NB	3.43	111.93	109.47
4	D	701	IMN	C7-C8-N	3.07	111.76	108.85
3	A	682	HEM	C3B-C4B-NB	2.97	111.60	109.47
2	C	671	NAG	C2-N2-C7	-2.89	119.03	122.90
4	D	701	IMN	C10-C9-N	2.80	123.30	117.92
4	B	701	IMN	C7-C8-N	2.80	111.51	108.85
4	C	701	IMN	C10-C9-N	2.73	123.16	117.92
4	A	701	IMN	C10-C9-N	2.72	123.14	117.92
4	B	701	IMN	C10-C9-N	2.63	122.97	117.92
4	A	701	IMN	C16-C8-C7	-2.63	123.96	128.07
4	C	701	IMN	C7-C8-N	2.61	111.33	108.85
2	D	671	NAG	C2-N2-C7	-2.59	119.42	122.90
4	A	701	IMN	C7-C8-N	2.59	111.31	108.85
4	C	701	IMN	C16-C8-C7	-2.59	124.02	128.07
2	C	671	NAG	C1-O5-C5	2.46	115.48	112.19
2	B	671	NAG	C2-N2-C7	-2.44	119.64	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	701	IMN	C16-C8-C7	-2.42	124.29	128.07
2	B	671	NAG	C1-O5-C5	2.38	115.38	112.19
4	D	701	IMN	C-C1-C7	-2.38	105.21	107.38
2	B	661	NAG	C2-N2-C7	-2.37	119.72	122.90
2	D	661	NAG	C2-N2-C7	-2.35	119.75	122.90
2	D	671	NAG	C1-O5-C5	2.29	115.25	112.19
4	D	701	IMN	C5-C-C1	-2.22	119.47	121.95
2	D	671	NAG	C6-C5-C4	-2.21	107.60	113.02
2	C	661	NAG	C1-O5-C5	-2.17	109.27	112.19
4	B	701	IMN	C-N-C8	2.12	109.53	107.80
4	C	701	IMN	C-C1-C7	-2.12	105.45	107.38
2	A	681	NAG	C4-C3-C2	2.11	114.10	111.02
2	A	661	NAG	O5-C1-C2	-2.09	108.06	111.29
4	B	701	IMN	C16-C8-C7	-2.07	124.83	128.07
3	B	682	HEM	CMD-C2D-C1D	2.07	128.26	125.03
4	A	701	IMN	C-C1-C7	-2.06	105.50	107.38
2	C	661	NAG	C2-N2-C7	-2.05	120.15	122.90
3	B	682	HEM	CAA-C2A-C1A	2.02	128.89	124.94
4	B	701	IMN	C-C1-C7	-2.02	105.54	107.38
4	C	701	IMN	C15-C14-C13	2.01	121.27	119.24

There are no chirality outliers.

All (68) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	701	IMN	C18-C17-C7-C1
2	D	661	NAG	O5-C5-C6-O6
2	B	671	NAG	O5-C5-C6-O6
2	C	661	NAG	O5-C5-C6-O6
2	B	661	NAG	O5-C5-C6-O6
2	C	671	NAG	O5-C5-C6-O6
2	A	661	NAG	O5-C5-C6-O6
2	A	671	NAG	O5-C5-C6-O6
2	D	671	NAG	O5-C5-C6-O6
2	B	671	NAG	C4-C5-C6-O6
2	C	671	NAG	C4-C5-C6-O6
2	A	671	NAG	C4-C5-C6-O6
2	D	671	NAG	C4-C5-C6-O6
4	A	701	IMN	C10-C9-N-C8
4	A	701	IMN	O1-C9-N-C8
4	B	701	IMN	C10-C9-N-C8
4	B	701	IMN	O1-C9-N-C8

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Mol	Chain	Res	Type	Atoms
4	C	701	IMN	C10-C9-N-C8
4	C	701	IMN	O1-C9-N-C8
4	D	701	IMN	C10-C9-N-C8
4	D	701	IMN	O1-C9-N-C8
4	A	701	IMN	C18-C17-C7-C1
2	D	661	NAG	C4-C5-C6-O6
2	B	661	NAG	C4-C5-C6-O6
4	A	701	IMN	C18-C17-C7-C8
4	B	701	IMN	C18-C17-C7-C8
4	C	701	IMN	C18-C17-C7-C8
4	D	701	IMN	C18-C17-C7-C8
2	C	661	NAG	C4-C5-C6-O6
2	A	661	NAG	C4-C5-C6-O6
4	B	701	IMN	C18-C17-C7-C1
4	D	701	IMN	C18-C17-C7-C1
3	A	682	HEM	C4C-C3C-CAC-CBC
4	A	701	IMN	C15-C10-C9-O1
4	A	701	IMN	O1-C9-N-C
4	B	701	IMN	O1-C9-N-C
4	C	701	IMN	O1-C9-N-C
4	D	701	IMN	O1-C9-N-C
4	C	701	IMN	C15-C10-C9-O1
4	B	701	IMN	C15-C10-C9-O1
4	D	701	IMN	C15-C10-C9-O1
4	A	701	IMN	C10-C9-N-C
4	C	701	IMN	C10-C9-N-C
4	D	701	IMN	C10-C9-N-C
2	A	681	NAG	C4-C5-C6-O6
4	A	701	IMN	C15-C10-C9-N
4	B	701	IMN	C15-C10-C9-N
4	C	701	IMN	C15-C10-C9-N
4	B	701	IMN	C10-C9-N-C
4	D	701	IMN	C15-C10-C9-N
4	A	701	IMN	C11-C10-C9-O1
2	A	681	NAG	O5-C5-C6-O6
3	B	682	HEM	C4C-C3C-CAC-CBC
3	C	682	HEM	CAD-CBD-CGD-O2D
3	C	682	HEM	CAA-CBA-CGA-O1A
3	D	682	HEM	CAD-CBD-CGD-O2D
4	D	701	IMN	C11-C10-C9-O1
4	C	701	IMN	C11-C10-C9-O1
4	B	701	IMN	C11-C10-C9-O1

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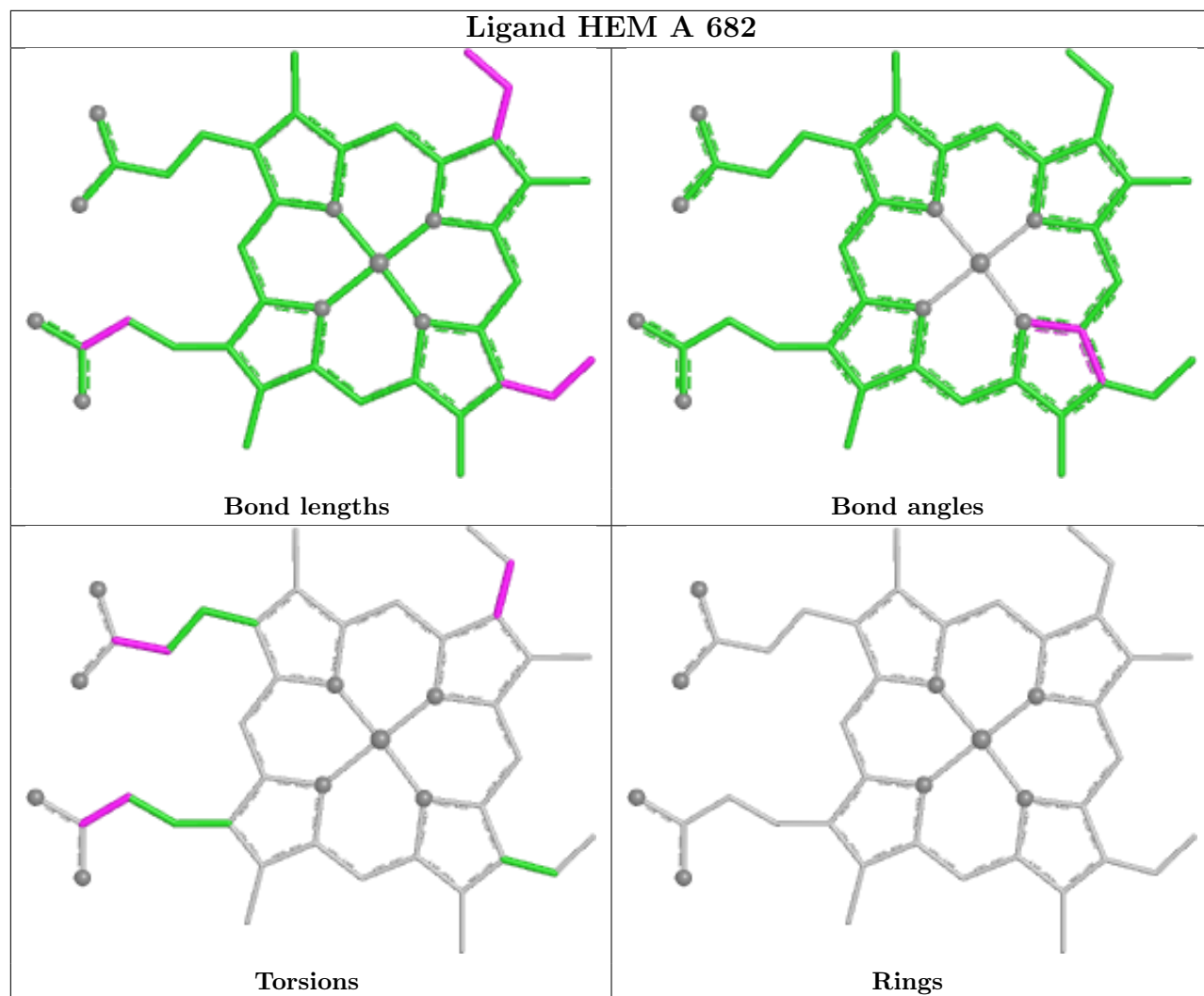
Mol	Chain	Res	Type	Atoms
3	A	682	HEM	CAA-CBA-CGA-O1A
3	C	682	HEM	CAA-CBA-CGA-O2A
3	A	682	HEM	CAD-CBD-CGD-O2D
3	B	682	HEM	CAD-CBD-CGD-O2D
3	C	682	HEM	CAD-CBD-CGD-O1D
3	B	682	HEM	CAA-CBA-CGA-O1A
3	D	682	HEM	CAA-CBA-CGA-O1A
3	A	682	HEM	CAA-CBA-CGA-O2A
4	A	701	IMN	C11-C10-C9-N

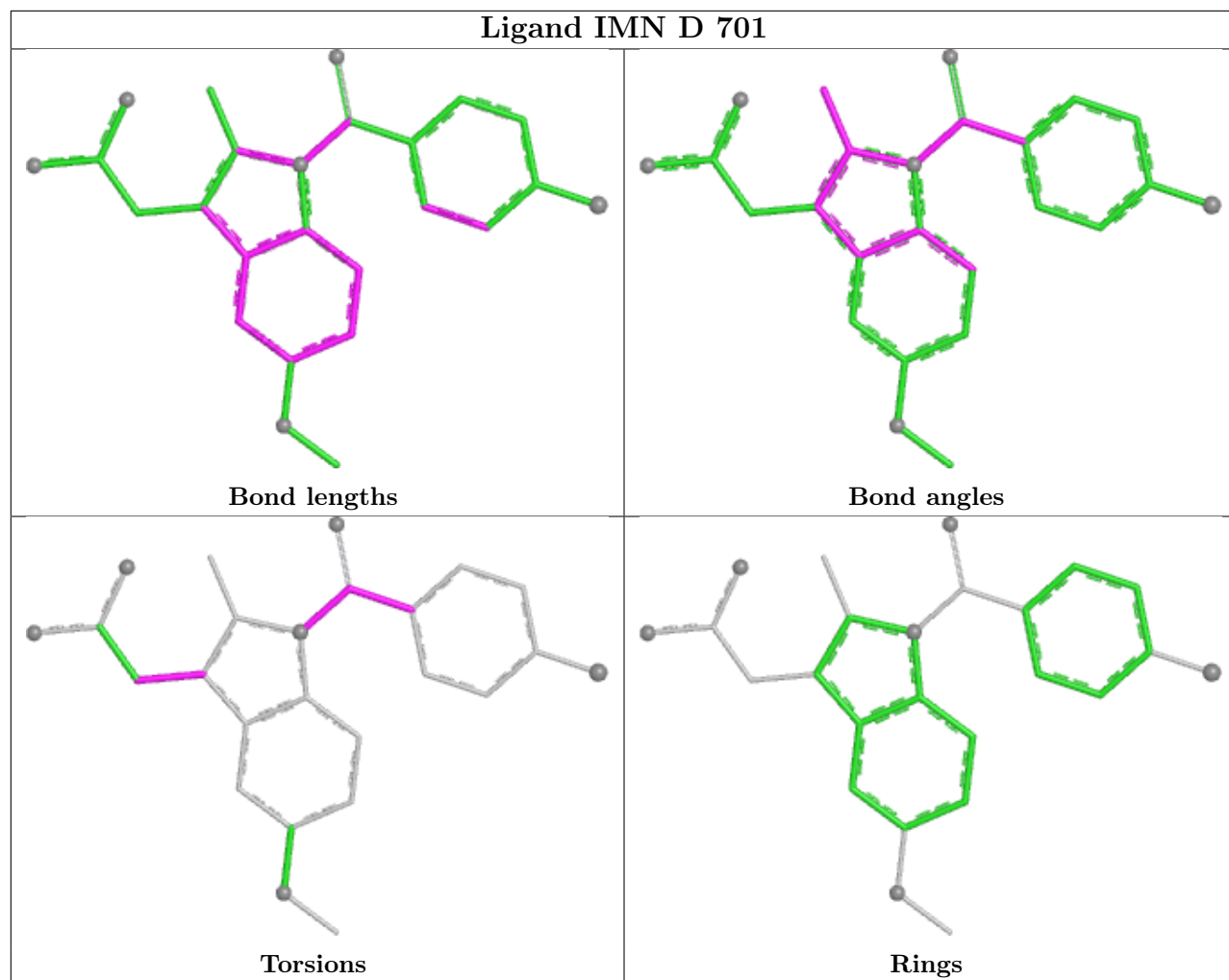
There are no ring outliers.

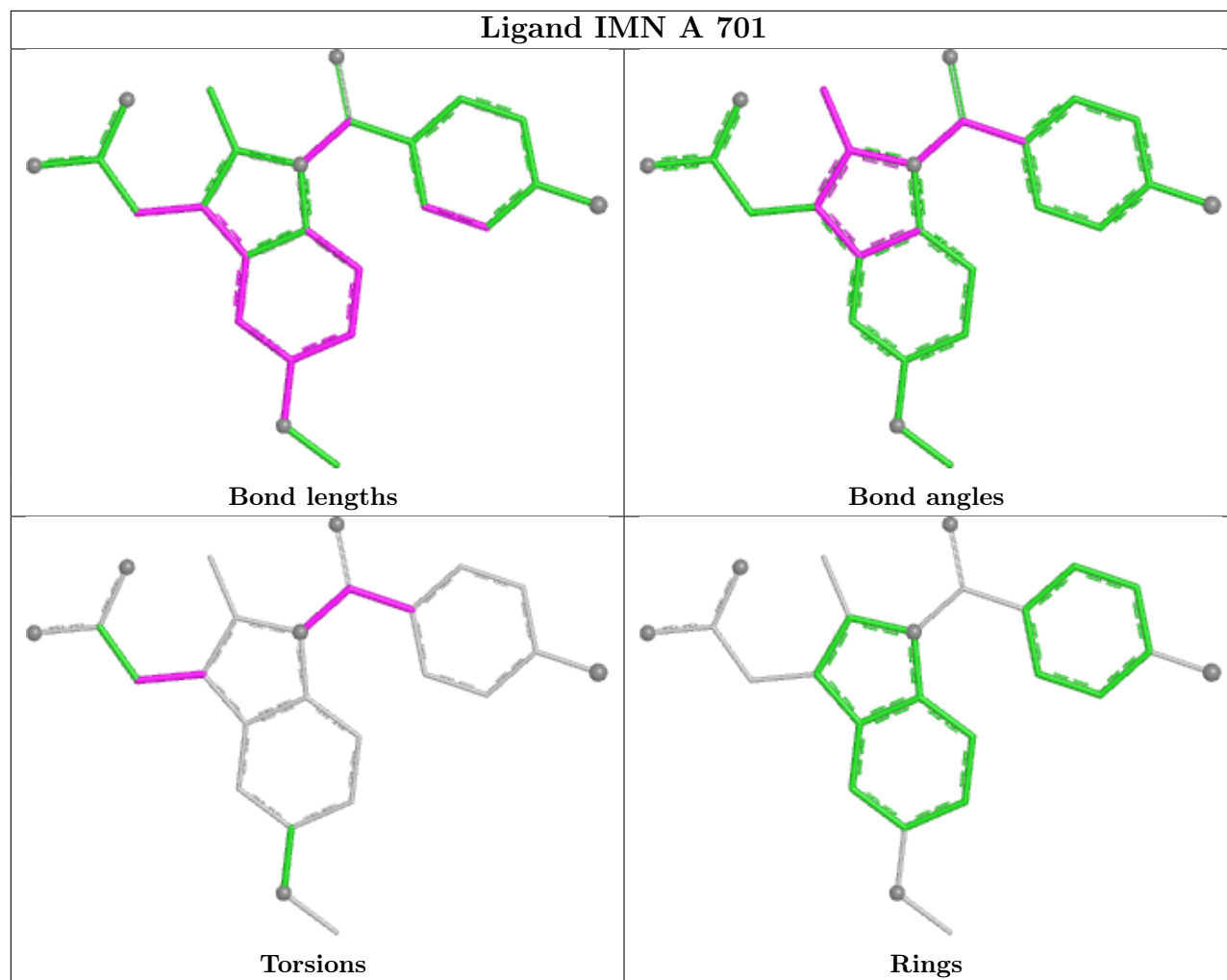
11 monomers are involved in 56 short contacts:

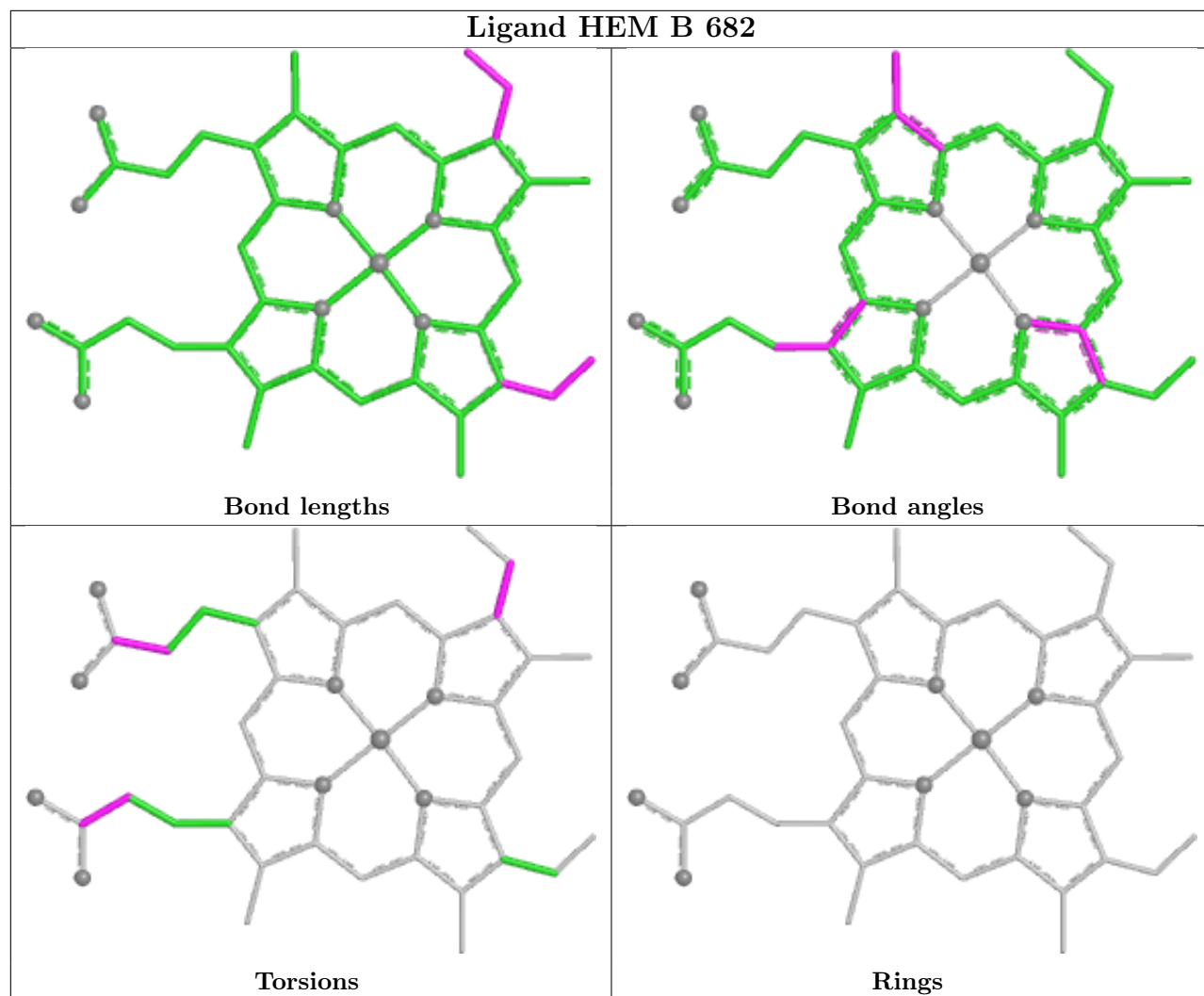
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	682	HEM	4	0
4	D	701	IMN	9	0
2	C	661	NAG	1	0
2	A	661	NAG	1	0
4	A	701	IMN	8	0
3	B	682	HEM	5	0
3	D	682	HEM	5	0
4	B	701	IMN	9	0
4	C	701	IMN	8	0
3	C	682	HEM	5	0
2	B	661	NAG	1	0

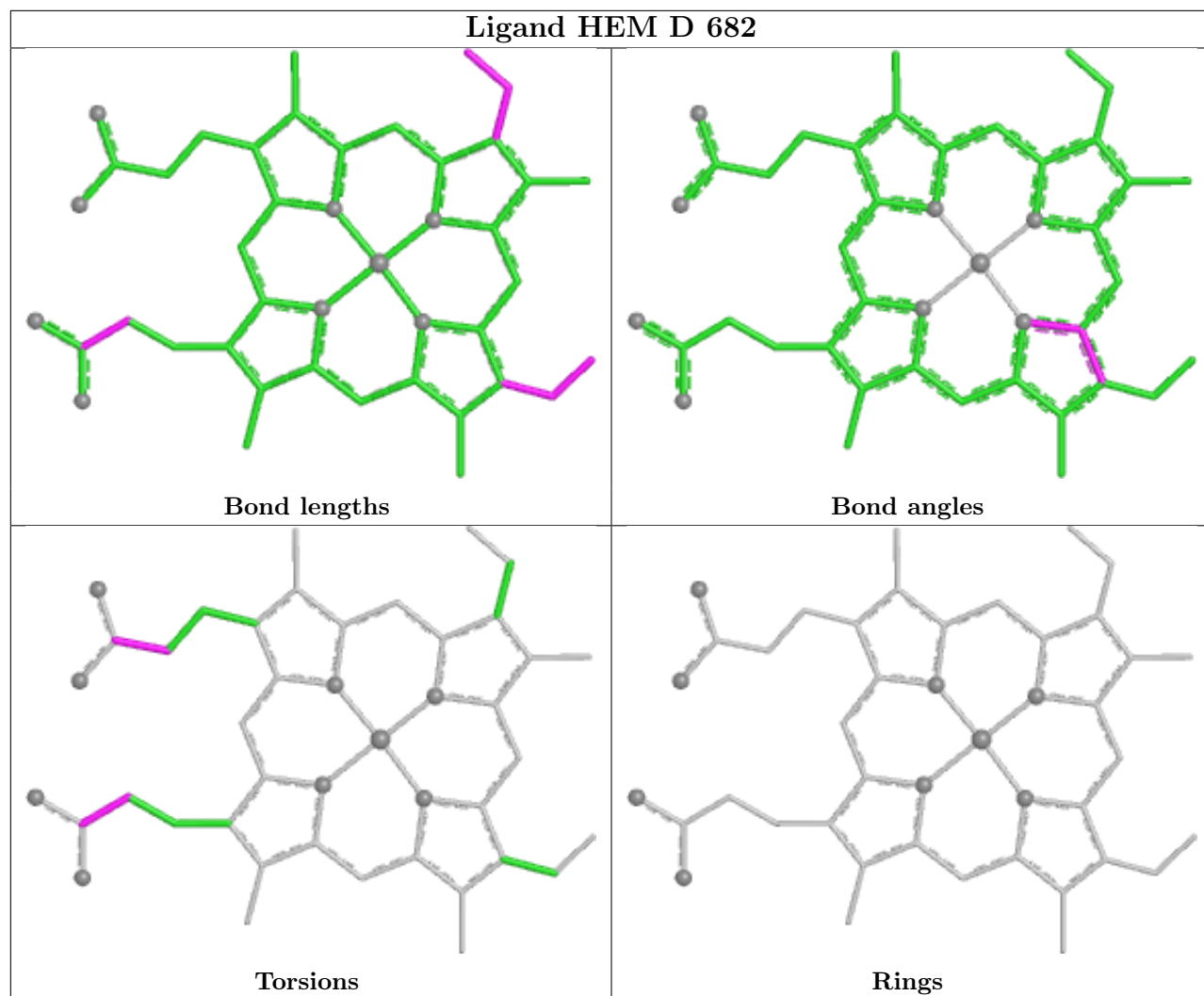
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

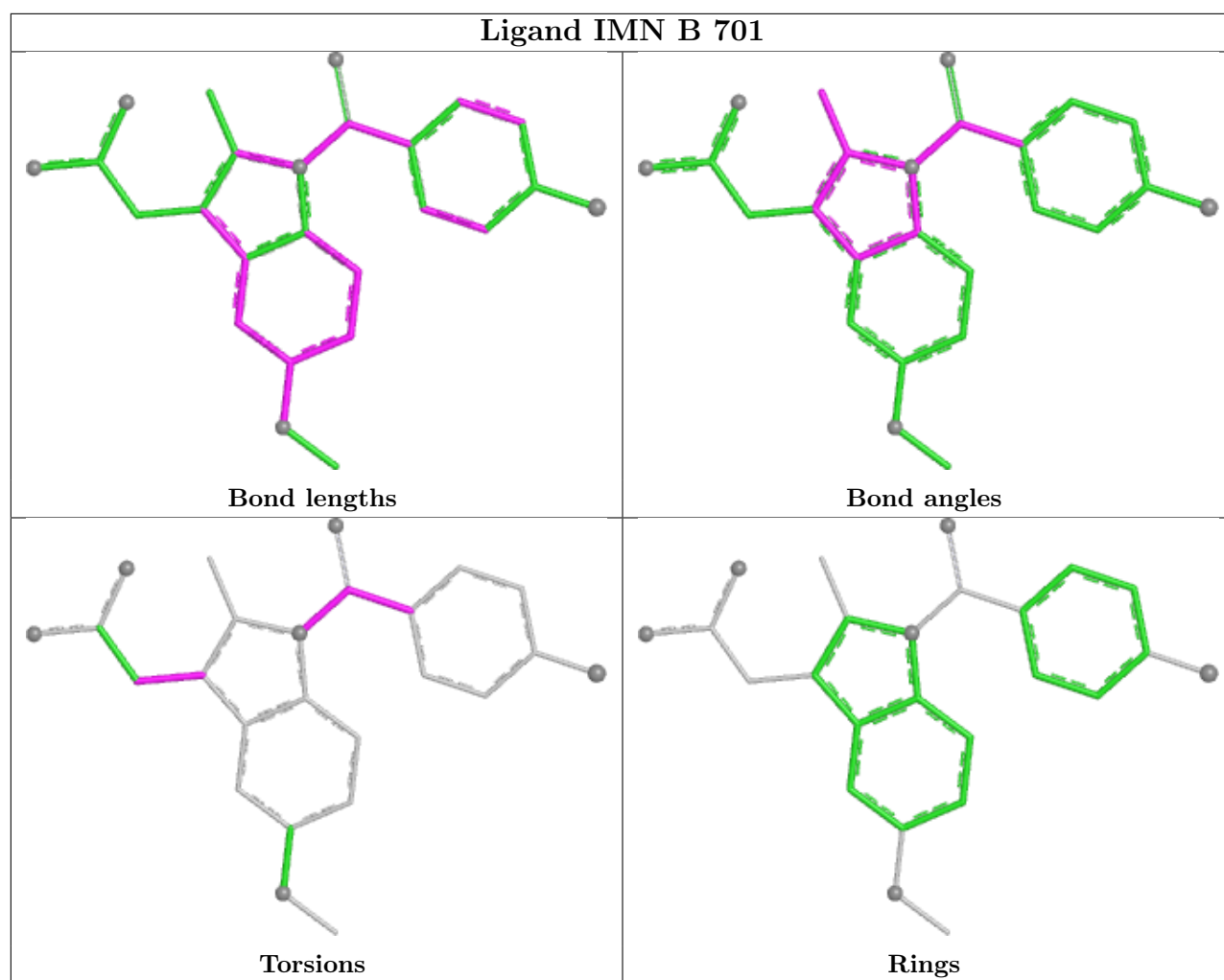


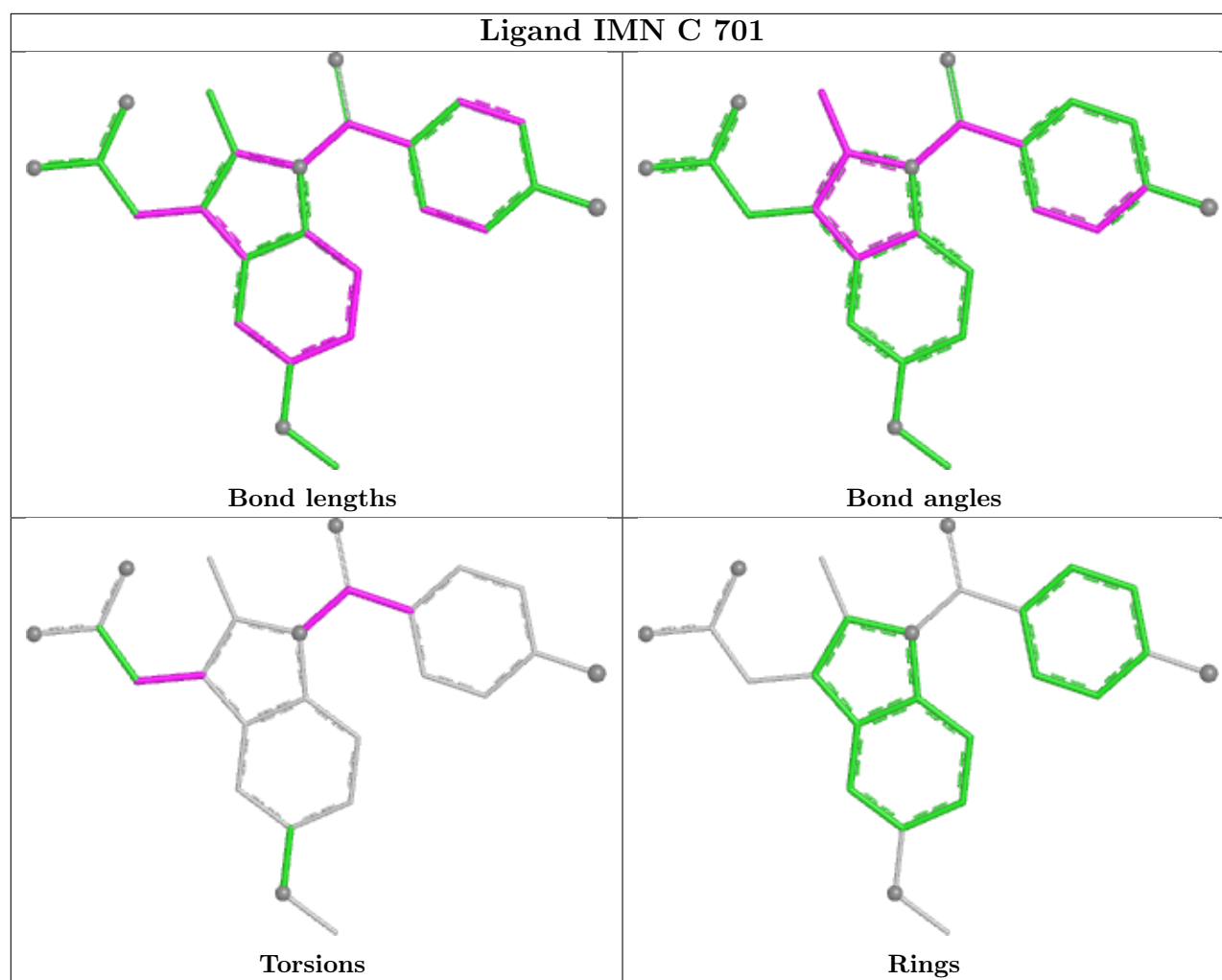


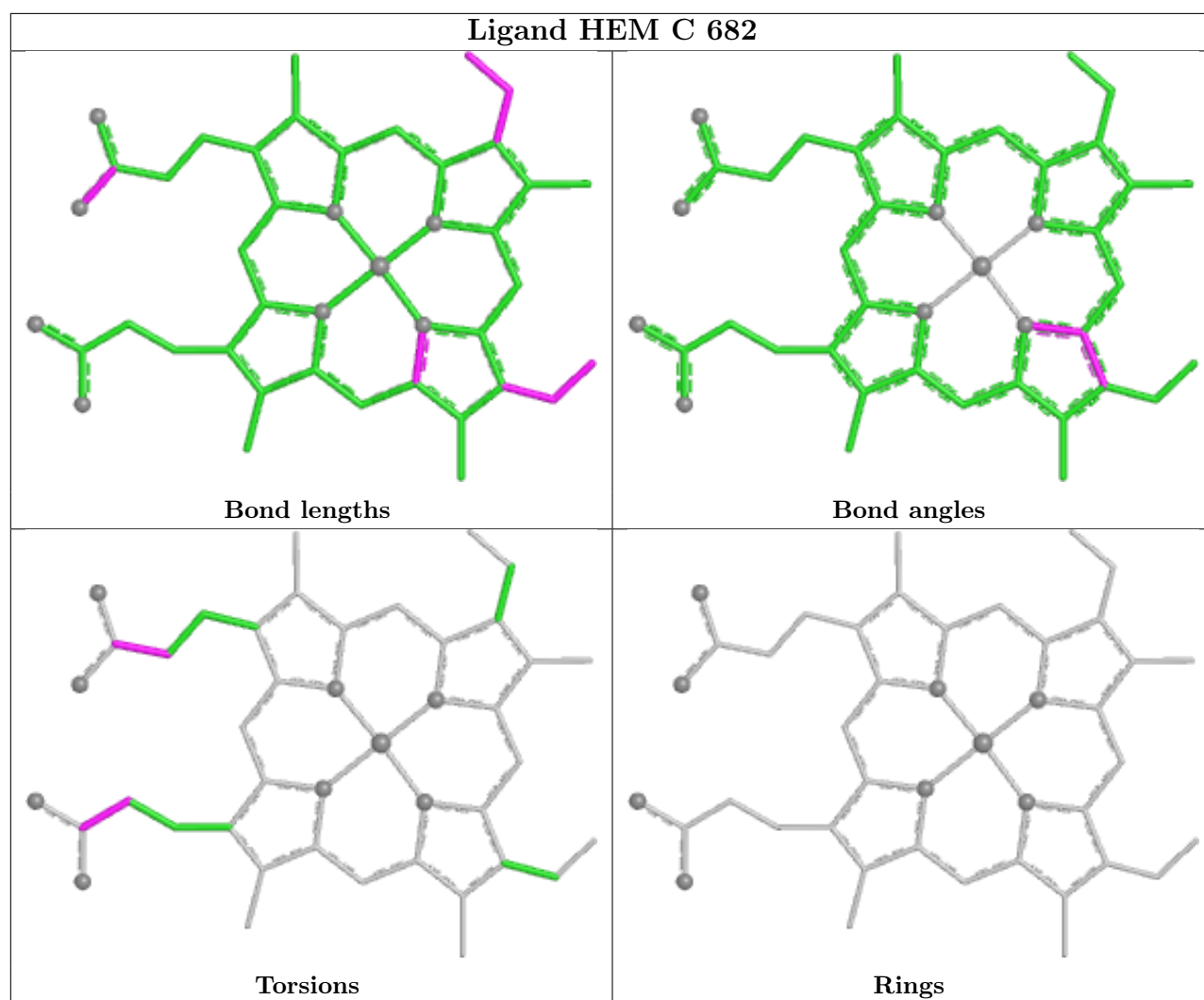












5.7 Other polymers [i](#)

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	552/587 (94%)	-0.07	4 (0%) 84 79	2, 10, 22, 35	0
1	B	552/587 (94%)	-0.02	2 (0%) 88 85	2, 10, 22, 34	0
1	C	552/587 (94%)	-0.07	5 (0%) 81 75	2, 10, 22, 34	0
1	D	552/587 (94%)	-0.08	4 (0%) 84 79	2, 10, 24, 31	0
All	All	2208/2348 (94%)	-0.06	15 (0%) 84 79	2, 10, 22, 35	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	450	ALA	2.8
1	A	400	GLN	2.6
1	D	543	GLN	2.5
1	C	161	THR	2.5
1	C	409	TYR	2.4
1	C	111	LEU	2.4
1	C	283	LEU	2.4
1	A	107	PHE	2.3
1	A	272	GLU	2.2
1	D	288	GLY	2.2
1	D	54	GLN	2.2
1	A	88	THR	2.1
1	D	398	GLU	2.1
1	B	450	ALA	2.0
1	B	442	ILE	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

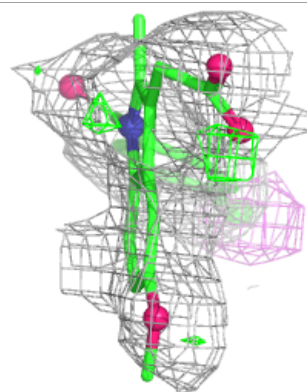
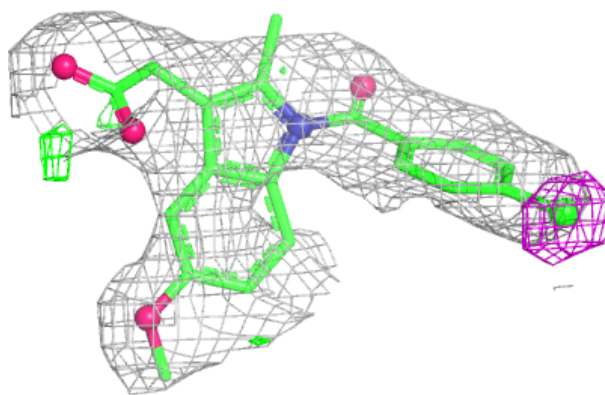
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	A	661	14/15	0.68	0.15	19,23,28,28	0
2	NAG	B	681	14/15	0.72	0.14	14,27,29,31	0
2	NAG	C	661	14/15	0.73	0.13	20,23,29,30	0
2	NAG	D	681	14/15	0.75	0.12	19,27,29,33	0
2	NAG	D	661	14/15	0.77	0.12	16,22,29,29	0
2	NAG	C	681	14/15	0.77	0.12	19,26,28,32	0
4	IMN	D	701	25/25	0.79	0.13	8,22,23,26	0
2	NAG	B	661	14/15	0.80	0.12	15,22,29,31	0
4	IMN	B	701	25/25	0.85	0.11	5,18,24,25	0
2	NAG	A	681	14/15	0.85	0.09	18,24,27,32	0
2	NAG	C	671	14/15	0.87	0.13	2,10,20,25	0
3	HEM	D	682	43/43	0.88	0.11	4,10,20,33	0
2	NAG	B	671	14/15	0.88	0.10	2,8,19,24	0
3	HEM	B	682	43/43	0.88	0.12	2,10,20,33	0
2	NAG	D	671	14/15	0.90	0.10	2,10,17,21	0
2	NAG	A	671	14/15	0.90	0.09	2,8,20,24	0
4	IMN	C	701	25/25	0.92	0.09	3,20,23,24	0
4	IMN	A	701	25/25	0.92	0.09	2,18,21,22	0
3	HEM	A	682	43/43	0.93	0.09	2,8,23,32	0
3	HEM	C	682	43/43	0.94	0.07	2,7,19,27	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

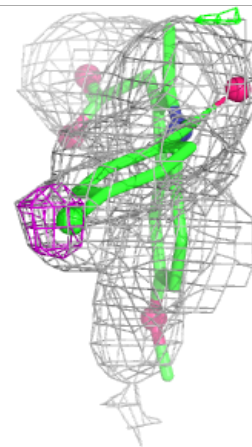
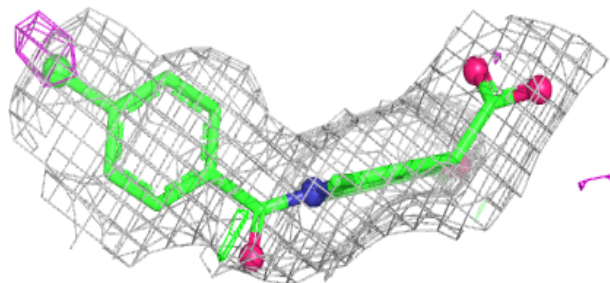
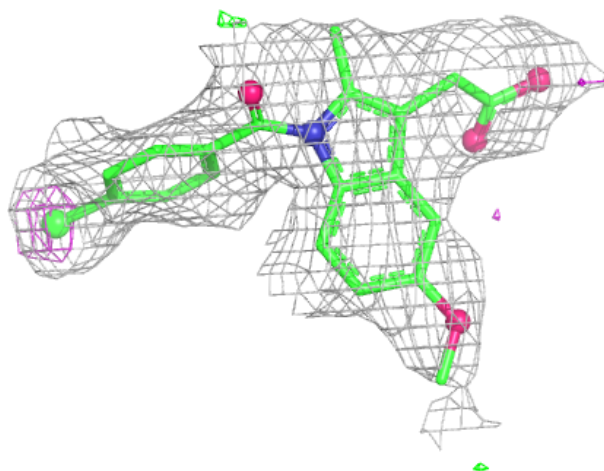
Electron density around IMN D 701:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



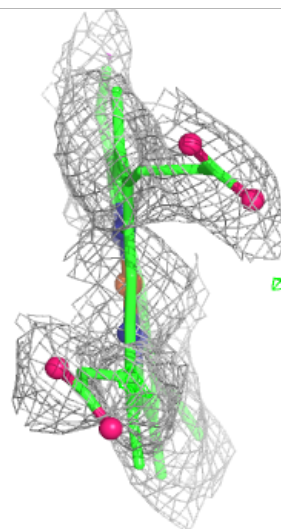
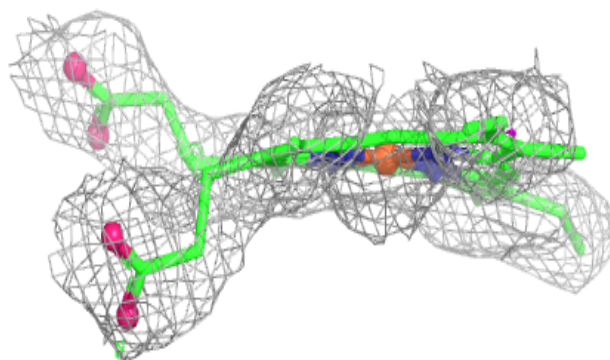
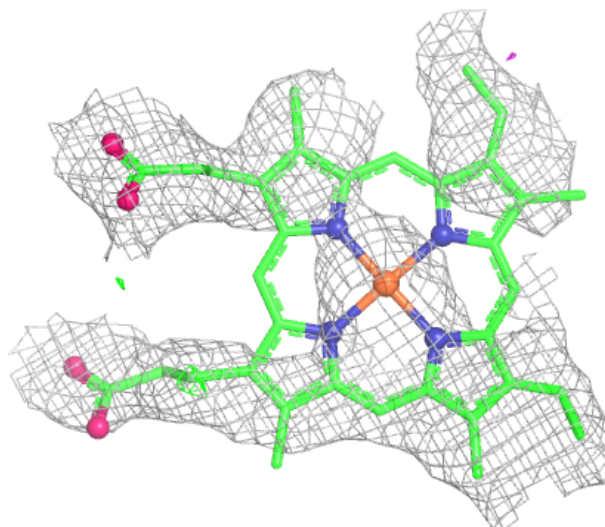
Electron density around IMN B 701:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



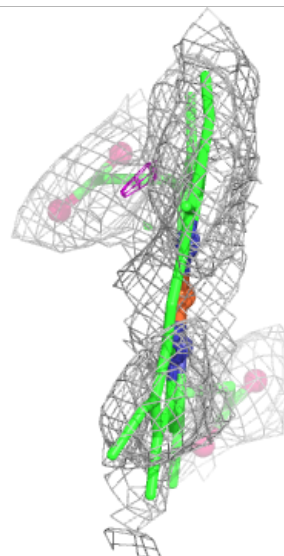
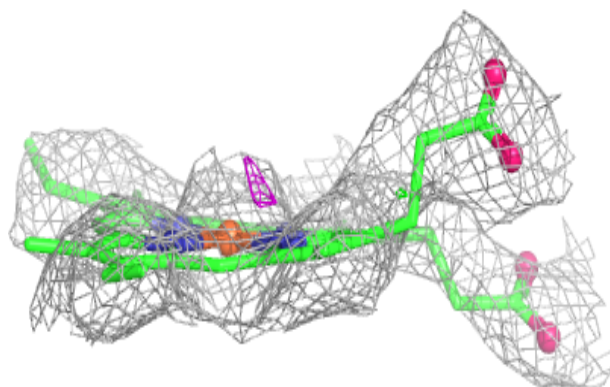
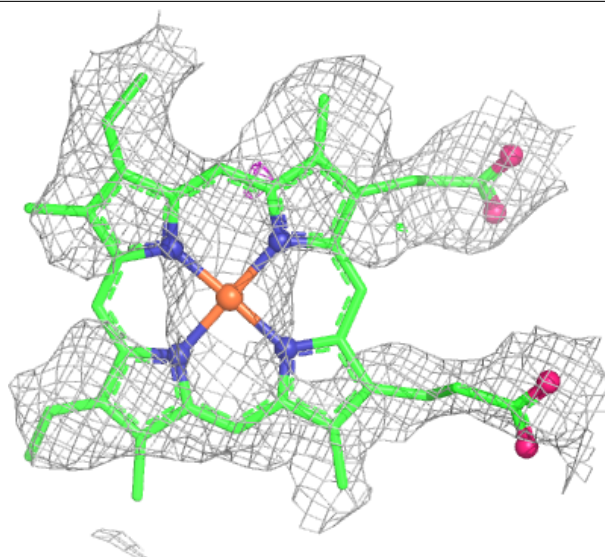
Electron density around HEM D 682:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



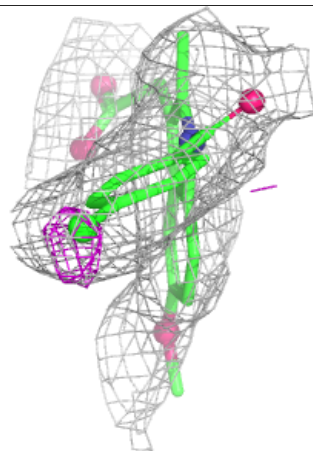
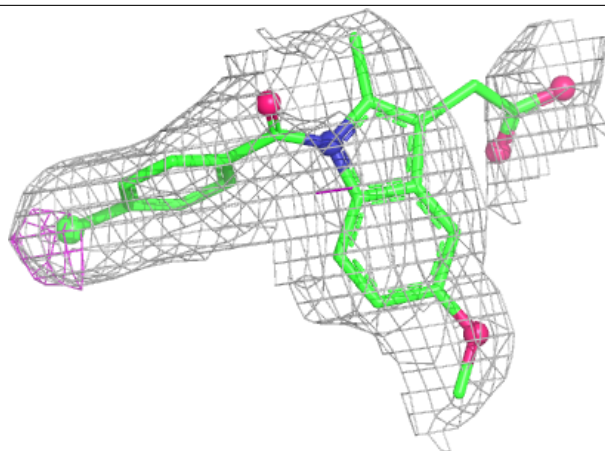
Electron density around HEM B 682:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

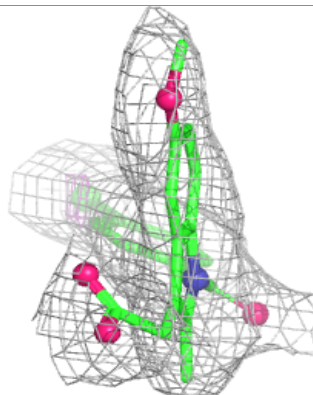
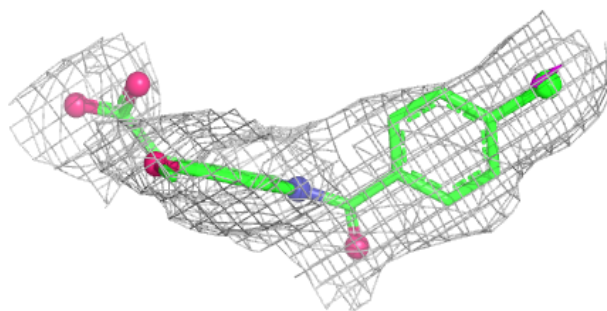
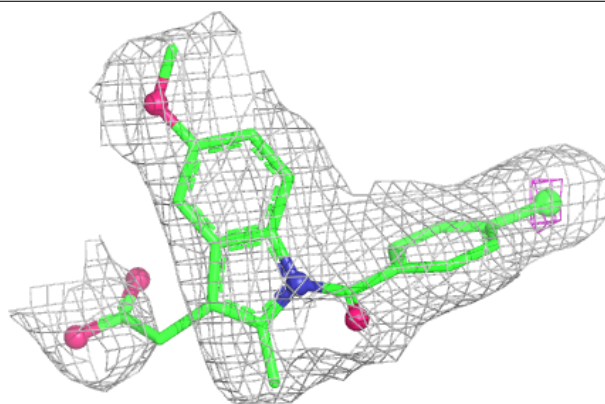


Electron density around IMN C 701:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

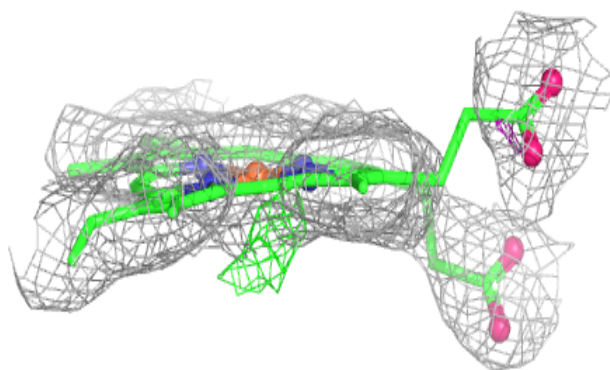
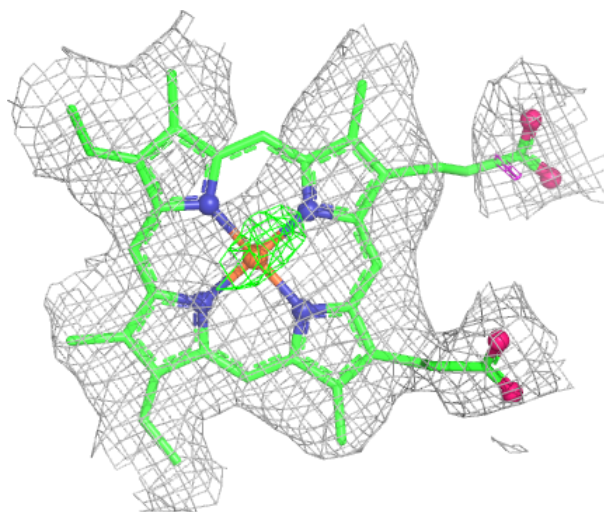
**Electron density around IMN A 701:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



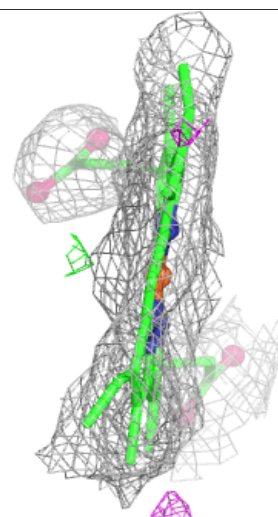
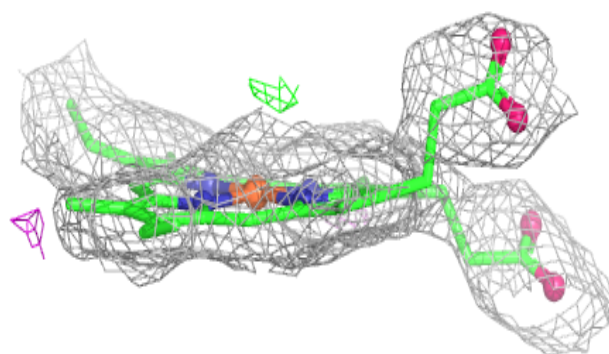
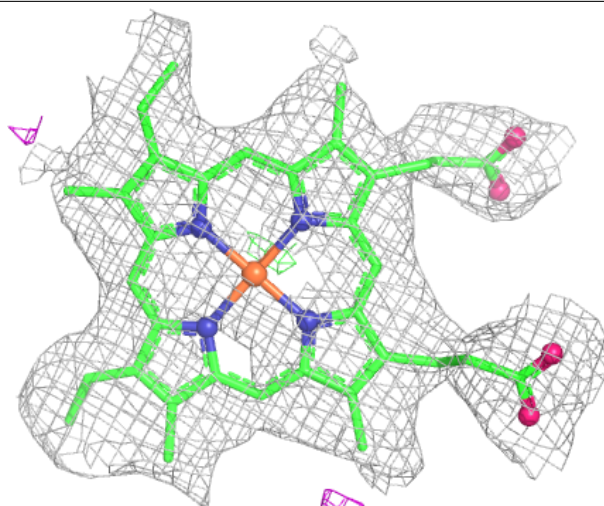
Electron density around HEM A 682:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM C 682:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.