



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:31 PM UTC

PDB ID : 9SI5 / pdb_00009si5
Title : Crystal structure of DoxA in complex with reaction intermediate DHD
Authors : Kim, R.Q.; Metsa-Ketela, M.
Deposited on : 2025-08-28
Resolution : 2.54 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

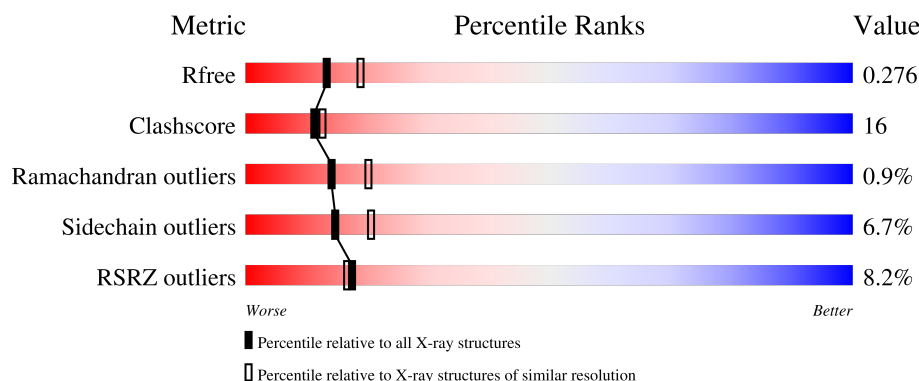
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	457	2% 60% 23% 5% 11%
1	D	457	12% 58% 25% • 11%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6585 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 Cytochrome P-450 monooxygenase DoxA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	407	Total	C	N	O	S	0	0	0
			3146	1991	573	571	11			
1	D	407	Total	C	N	O	S	0	0	0
			3141	1988	572	570	11			

There are 86 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-34	MET	-	initiating methionine	UNP Q9ZAU3
A	-33	GLY	-	expression tag	UNP Q9ZAU3
A	-32	GLY	-	expression tag	UNP Q9ZAU3
A	-31	SER	-	expression tag	UNP Q9ZAU3
A	-30	HIS	-	expression tag	UNP Q9ZAU3
A	-29	HIS	-	expression tag	UNP Q9ZAU3
A	-28	HIS	-	expression tag	UNP Q9ZAU3
A	-27	HIS	-	expression tag	UNP Q9ZAU3
A	-26	HIS	-	expression tag	UNP Q9ZAU3
A	-25	HIS	-	expression tag	UNP Q9ZAU3
A	-24	GLY	-	expression tag	UNP Q9ZAU3
A	-23	MET	-	expression tag	UNP Q9ZAU3
A	-22	ALA	-	expression tag	UNP Q9ZAU3
A	-21	SER	-	expression tag	UNP Q9ZAU3
A	-20	MET	-	expression tag	UNP Q9ZAU3
A	-19	THR	-	expression tag	UNP Q9ZAU3
A	-18	GLY	-	expression tag	UNP Q9ZAU3
A	-17	GLY	-	expression tag	UNP Q9ZAU3
A	-16	GLN	-	expression tag	UNP Q9ZAU3
A	-15	GLN	-	expression tag	UNP Q9ZAU3
A	-14	MET	-	expression tag	UNP Q9ZAU3
A	-13	GLY	-	expression tag	UNP Q9ZAU3
A	-12	ARG	-	expression tag	UNP Q9ZAU3
A	-11	ASP	-	expression tag	UNP Q9ZAU3
A	-10	LEU	-	expression tag	UNP Q9ZAU3

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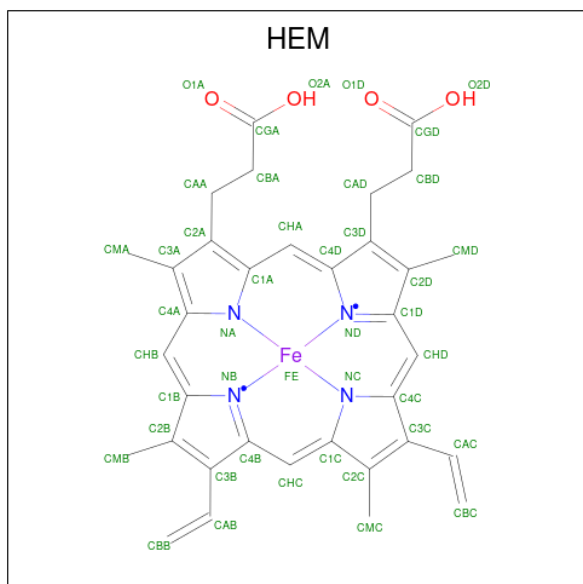
Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	TYR	-	expression tag	UNP Q9ZAU3
A	-8	ASP	-	expression tag	UNP Q9ZAU3
A	-7	ASP	-	expression tag	UNP Q9ZAU3
A	-6	ASP	-	expression tag	UNP Q9ZAU3
A	-5	ASP	-	expression tag	UNP Q9ZAU3
A	-4	LYS	-	expression tag	UNP Q9ZAU3
A	-3	ASP	-	expression tag	UNP Q9ZAU3
A	-2	PRO	-	expression tag	UNP Q9ZAU3
A	-1	SER	-	expression tag	UNP Q9ZAU3
A	0	SER	-	expression tag	UNP Q9ZAU3
A	1	ARG	-	expression tag	UNP Q9ZAU3
A	2	SER	-	expression tag	UNP Q9ZAU3
A	3	GLY	-	expression tag	UNP Q9ZAU3
A	4	GLU	-	expression tag	UNP Q9ZAU3
A	5	ALA	-	expression tag	UNP Q9ZAU3
A	6	PRO	-	expression tag	UNP Q9ZAU3
A	7	ARG	-	expression tag	UNP Q9ZAU3
A	8	VAL	-	expression tag	UNP Q9ZAU3
D	-34	MET	-	initiating methionine	UNP Q9ZAU3
D	-33	GLY	-	expression tag	UNP Q9ZAU3
D	-32	GLY	-	expression tag	UNP Q9ZAU3
D	-31	SER	-	expression tag	UNP Q9ZAU3
D	-30	HIS	-	expression tag	UNP Q9ZAU3
D	-29	HIS	-	expression tag	UNP Q9ZAU3
D	-28	HIS	-	expression tag	UNP Q9ZAU3
D	-27	HIS	-	expression tag	UNP Q9ZAU3
D	-26	HIS	-	expression tag	UNP Q9ZAU3
D	-25	HIS	-	expression tag	UNP Q9ZAU3
D	-24	GLY	-	expression tag	UNP Q9ZAU3
D	-23	MET	-	expression tag	UNP Q9ZAU3
D	-22	ALA	-	expression tag	UNP Q9ZAU3
D	-21	SER	-	expression tag	UNP Q9ZAU3
D	-20	MET	-	expression tag	UNP Q9ZAU3
D	-19	THR	-	expression tag	UNP Q9ZAU3
D	-18	GLY	-	expression tag	UNP Q9ZAU3
D	-17	GLY	-	expression tag	UNP Q9ZAU3
D	-16	GLN	-	expression tag	UNP Q9ZAU3
D	-15	GLN	-	expression tag	UNP Q9ZAU3
D	-14	MET	-	expression tag	UNP Q9ZAU3
D	-13	GLY	-	expression tag	UNP Q9ZAU3
D	-12	ARG	-	expression tag	UNP Q9ZAU3
D	-11	ASP	-	expression tag	UNP Q9ZAU3

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-10	LEU	-	expression tag	UNP Q9ZAU3
D	-9	TYR	-	expression tag	UNP Q9ZAU3
D	-8	ASP	-	expression tag	UNP Q9ZAU3
D	-7	ASP	-	expression tag	UNP Q9ZAU3
D	-6	ASP	-	expression tag	UNP Q9ZAU3
D	-5	ASP	-	expression tag	UNP Q9ZAU3
D	-4	LYS	-	expression tag	UNP Q9ZAU3
D	-3	ASP	-	expression tag	UNP Q9ZAU3
D	-2	PRO	-	expression tag	UNP Q9ZAU3
D	-1	SER	-	expression tag	UNP Q9ZAU3
D	0	SER	-	expression tag	UNP Q9ZAU3
D	1	ARG	-	expression tag	UNP Q9ZAU3
D	2	SER	-	expression tag	UNP Q9ZAU3
D	3	GLY	-	expression tag	UNP Q9ZAU3
D	4	GLU	-	expression tag	UNP Q9ZAU3
D	5	ALA	-	expression tag	UNP Q9ZAU3
D	6	PRO	-	expression tag	UNP Q9ZAU3
D	7	ARG	-	expression tag	UNP Q9ZAU3
D	8	VAL	-	expression tag	UNP Q9ZAU3

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$) (labeled as "Ligand of Interest" by depositor).



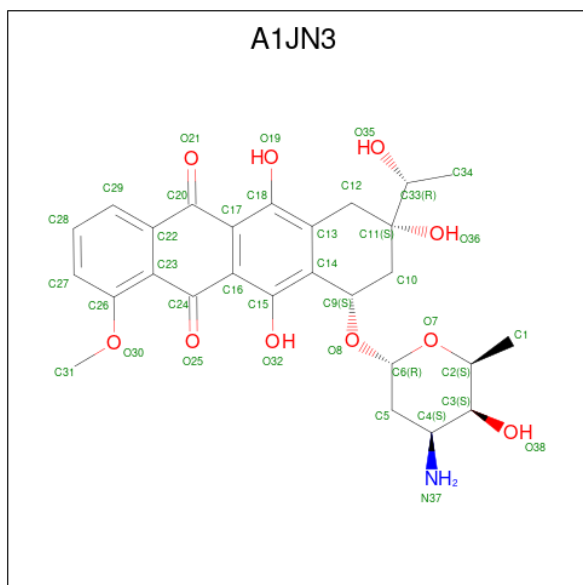
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	
			43	34	1	4	4	
							0	0

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

- Molecule 3 is (7 {S},9 {S})-7-[(2 {R},4 {S},5 {S},6 {S})-4-azanyl-6-methyl-5-oxidanyl-oxan-2-yl]oxy-4-methoxy-6,9,11-tris(oxidanyl)-9-[(1 {R})-1-oxidanylethyl]-8,10-dihydro-7 {H}-tetracene-5,12-dione (CCD ID: A1JN3) (formula: C₂₇H₃₁NO₁₀) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			38	27	1	10		
3	D	1	Total	C	N	O	0	0
			38	27	1	10		

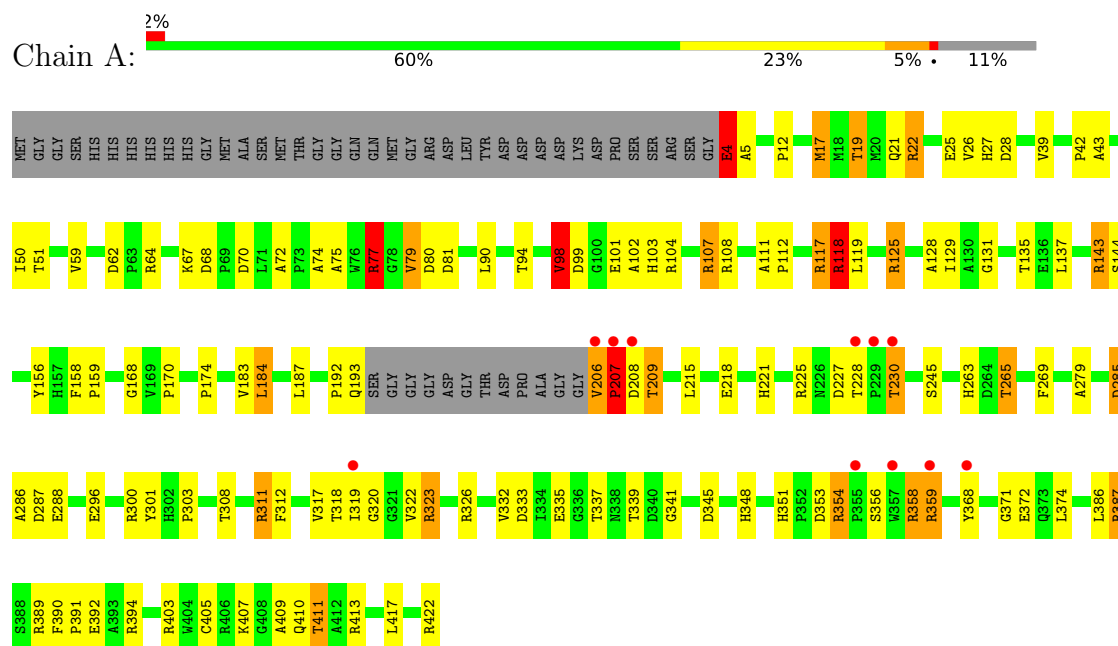
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	87	Total	O	0	0
			87	87		
4	D	49	Total	O	0	0
			49	49		

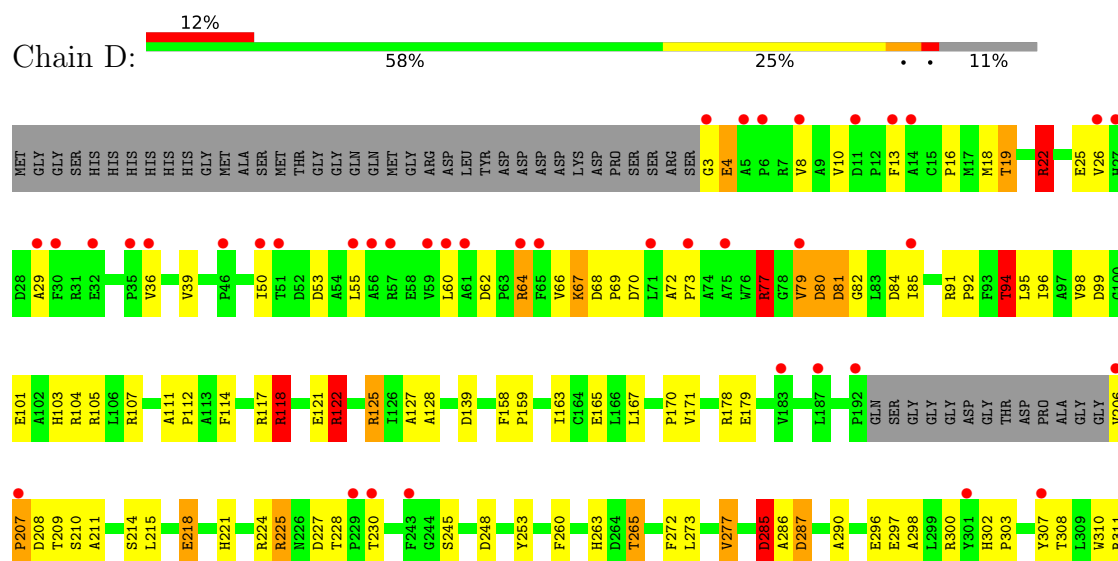
3 Residue-property plots [i](#)

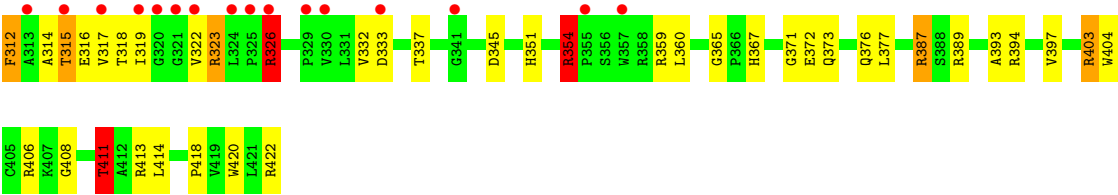
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Cytochrome P-450 monooxygenase DoxA



• Molecule 1: Cytochrome P-450 monooxygenase DoxA





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	101.87Å 108.06Å 179.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	89.89 – 2.54 89.89 – 2.54	Depositor EDS
% Data completeness (in resolution range)	98.9 (89.89-2.54) 98.1 (89.89-2.54)	Depositor EDS
R_{merge}	0.30	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.08 (at 2.55Å)	Xtriage
Refinement program	REFMAC 5.8.0430 (refmacat 0.4.105)	Depositor
R, R_{free}	0.234 , 0.281 0.232 , 0.276	Depositor DCC
R_{free} test set	1648 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	51.2	Xtriage
Anisotropy	0.919	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 70.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6585	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.83% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: A1JN3, HEM

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	0.87	0/3225	1.59	42/4397 (1.0%)
1	D	0.77	1/3220 (0.0%)	1.48	35/4390 (0.8%)
All	All	0.82	1/6445 (0.0%)	1.54	77/8787 (0.9%)

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	10
1	D	0	12
All	All	0	22

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	118	ARG	NE-CZ	-5.03	1.27	1.33

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	118	ARG	CD-NE-CZ	-14.60	103.97	124.40
1	A	118	ARG	CD-NE-CZ	-10.58	109.58	124.40
1	A	22	ARG	NE-CZ-NH2	10.37	128.54	119.20
1	A	354	ARG	N-CA-CB	-10.19	94.64	110.03
1	D	22	ARG	NE-CZ-NH2	9.68	127.91	119.20
1	A	98	VAL	N-CA-CB	-9.13	95.46	112.36
1	A	265	THR	OG1-CB-CG2	-9.10	91.09	109.30
1	A	209	THR	CA-CB-OG1	-9.02	96.07	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	228	THR	CA-CB-OG1	-8.66	96.61	109.60
1	D	4	GLU	CB-CA-C	8.46	127.64	111.48
1	A	118	ARG	CA-CB-CG	-8.44	97.22	114.10
1	D	218	GLU	CB-CA-C	-8.41	96.84	110.79
1	D	285	ASP	CB-CA-C	-8.40	96.78	111.22
1	A	118	ARG	CB-CG-CD	8.24	130.26	111.30
1	D	114	PHE	CA-CB-CG	-7.92	105.88	113.80
1	D	22	ARG	NE-CZ-NH1	-7.73	113.77	121.50
1	D	228	THR	CA-CB-OG1	-7.68	98.08	109.60
1	D	94	THR	OG1-CB-CG2	-7.64	94.02	109.30
1	D	139	ASP	CA-CB-CG	7.53	120.13	112.60
1	D	326	ARG	CG-CD-NE	-6.89	96.83	112.00
1	D	315	THR	OG1-CB-CG2	-6.75	95.81	109.30
1	D	411	THR	OG1-CB-CG2	-6.74	95.82	109.30
1	D	209	THR	CA-CB-OG1	-6.69	99.57	109.60
1	A	192	PRO	N-CA-C	6.68	121.74	111.11
1	A	28	ASP	CA-CB-CG	6.66	119.25	112.60
1	A	288	GLU	N-CA-CB	-6.50	100.26	109.94
1	A	387	ARG	CB-CA-C	6.47	122.69	109.55
1	A	174	PRO	CB-CA-C	6.45	121.53	112.11
1	D	19	THR	OG1-CB-CG2	-6.38	96.54	109.30
1	D	218	GLU	CG-CD-OE1	-6.31	103.89	118.40
1	A	62	ASP	CA-CB-CG	6.29	118.89	112.60
1	D	387	ARG	NE-CZ-NH1	-6.25	115.25	121.50
1	D	117	ARG	N-CA-CB	6.24	119.40	110.16
1	A	387	ARG	NE-CZ-NH1	-6.21	115.29	121.50
1	A	337	THR	CA-CB-OG1	6.12	118.78	109.60
1	A	269	PHE	CA-CB-CG	-6.11	107.69	113.80
1	D	260	PHE	N-CA-CB	5.98	118.74	110.07
1	A	353	ASP	CA-CB-CG	5.95	118.55	112.60
1	A	101	GLU	CB-CG-CD	5.92	122.66	112.60
1	A	411	THR	OG1-CB-CG2	-5.90	97.50	109.30
1	D	118	ARG	NE-CZ-NH2	-5.87	113.92	119.20
1	D	170	PRO	CB-CA-C	5.85	116.73	111.40
1	D	80	ASP	CA-CB-CG	5.83	118.44	112.60
1	D	121	GLU	CB-CG-CD	5.74	122.35	112.60
1	D	4	GLU	N-CA-CB	-5.70	101.94	110.60
1	D	122	ARG	CA-CB-CG	-5.66	102.78	114.10
1	A	4	GLU	N-CA-CB	-5.65	100.90	110.50
1	A	104	ARG	CB-CA-C	5.64	120.15	110.79
1	A	207	PRO	CB-CA-C	5.63	120.86	111.56
1	D	62	ASP	CA-CB-CG	5.58	118.18	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	101	GLU	CB-CG-CD	5.53	122.01	112.60
1	A	323	ARG	CD-NE-CZ	-5.52	116.67	124.40
1	A	279	ALA	N-CA-C	-5.48	106.60	113.28
1	A	4	GLU	CA-CB-CG	-5.38	103.35	114.10
1	A	417	LEU	N-CA-CB	-5.37	104.42	111.23
1	A	359	ARG	CG-CD-NE	-5.36	100.20	112.00
1	D	394	ARG	NE-CZ-NH2	5.36	124.02	119.20
1	D	178	ARG	CG-CD-NE	-5.35	100.23	112.00
1	D	19	THR	CB-CA-C	5.35	118.62	109.80
1	A	137	LEU	N-CA-CB	-5.34	102.28	110.12
1	A	143	ARG	CB-CA-C	-5.33	101.94	110.79
1	A	337	THR	OG1-CB-CG2	-5.32	98.66	109.30
1	D	287	ASP	CA-CB-CG	5.32	117.92	112.60
1	A	94	THR	CA-CB-OG1	-5.31	101.64	109.60
1	A	407	LYS	N-CA-CB	5.26	119.20	110.47
1	A	99	ASP	CA-CB-CG	5.26	117.86	112.60
1	D	277	VAL	N-CA-CB	5.24	116.68	110.55
1	D	296	GLU	CB-CA-C	5.20	119.43	110.79
1	D	272	PHE	N-CA-CB	5.20	117.76	110.12
1	D	354	ARG	CB-CA-C	-5.19	99.95	110.17
1	A	184	LEU	N-CA-C	-5.18	105.54	111.14
1	A	19	THR	CA-CB-OG1	-5.17	101.84	109.60
1	A	51	THR	CA-CB-OG1	-5.17	101.85	109.60
1	A	296	GLU	CB-CA-C	5.16	119.31	110.74
1	A	230	THR	OG1-CB-CG2	-5.16	98.99	109.30
1	A	125	ARG	N-CA-CB	5.05	117.55	110.12
1	A	117	ARG	N-CA-CB	5.04	117.32	110.01

There are no chirality outliers.

All (22) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	107	ARG	Sidechain
1	A	118	ARG	Sidechain
1	A	311	ARG	Sidechain
1	A	323	ARG	Sidechain
1	A	358	ARG	Sidechain
1	A	359	ARG	Sidechain
1	A	394	ARG	Sidechain
1	A	422	ARG	Sidechain
1	A	64	ARG	Sidechain
1	A	77	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	D	118	ARG	Sidechain
1	D	122	ARG	Sidechain
1	D	125	ARG	Sidechain
1	D	22	ARG	Sidechain
1	D	225	ARG	Sidechain
1	D	323	ARG	Sidechain
1	D	326	ARG	Sidechain
1	D	359	ARG	Sidechain
1	D	403	ARG	Sidechain
1	D	422	ARG	Sidechain
1	D	64	ARG	Sidechain
1	D	77	ARG	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	3146	0	3120	78	0
1	D	3141	0	3115	120	1
2	A	43	0	30	4	0
2	D	43	0	30	7	0
3	A	38	0	0	3	0
3	D	38	0	0	4	0
4	A	87	0	0	8	0
4	D	49	0	0	6	0
All	All	6585	0	6295	206	1

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 16.

All (206) 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:D:206:VAL:O	1:D:208:ASP:N	1.64	1.27
1:A:206:VAL:O	1:A:208:ASP:N	1.84	1.11
1:D:80:ASP:O	1:D:82:GLY:N	1.89	1.05

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:206:VAL:C	1:D:208:ASP:H	1.64	1.02
1:D:206:VAL:N	1:D:207:PRO:HD2	1.74	1.01
1:D:3:GLY:N	1:D:4:GLU:OE2	1.98	0.96
1:D:4:GLU:N	1:D:4:GLU:CD	2.24	0.95
1:D:3:GLY:C	1:D:4:GLU:CD	2.35	0.94
1:D:206:VAL:N	1:D:207:PRO:CD	2.31	0.94
1:D:64:ARG:CZ	1:D:317:VAL:HG21	1.99	0.91
1:A:387:ARG:NH1	4:A:602:HOH:O	2.06	0.87
1:D:211:ALA:O	1:D:215:LEU:HD23	1.74	0.86
2:D:501:HEM:HMC1	2:D:501:HEM:HBC2	1.58	0.86
1:D:73:PRO:HD3	4:D:611:HOH:O	1.76	0.86
1:A:144:SER:O	4:A:601:HOH:O	1.95	0.84
1:D:367:HIS:HD2	2:D:501:HEM:O1D	1.61	0.84
1:D:72:ALA:HA	1:D:312:PHE:HE2	1.45	0.81
2:D:501:HEM:HBB2	2:D:501:HEM:HMB2	1.63	0.81
1:A:206:VAL:C	1:A:208:ASP:H	1.84	0.81
1:A:308:THR:HG23	1:A:332:VAL:HB	1.63	0.80
1:D:4:GLU:N	1:D:4:GLU:OE1	2.18	0.75
1:D:80:ASP:O	1:D:80:ASP:OD1	2.04	0.75
1:A:206:VAL:O	1:A:206:VAL:CG2	2.36	0.72
1:D:367:HIS:CD2	2:D:501:HEM:O1D	2.43	0.72
1:D:72:ALA:HA	1:D:312:PHE:CE2	2.25	0.72
1:D:297:GLU:OE2	1:D:354:ARG:HD2	1.90	0.70
1:A:4:GLU:HA	1:A:4:GLU:OE1	1.93	0.69
1:D:354:ARG:HH11	1:D:354:ARG:HG2	1.59	0.68
1:D:13:PHE:HB3	1:D:408:GLY:HA2	1.75	0.68
1:D:66:VAL:HG22	1:D:68:ASP:H	1.61	0.66
1:A:285:ASP:N	1:A:285:ASP:OD1	2.28	0.66
1:A:183:VAL:HG21	1:A:207:PRO:HG2	1.77	0.65
1:D:314:ALA:O	1:D:326:ARG:NH2	2.29	0.65
1:A:308:THR:CG2	1:A:332:VAL:HB	2.25	0.64
1:D:206:VAL:C	1:D:208:ASP:N	2.32	0.64
2:D:501:HEM:HBC2	2:D:501:HEM:CMC	2.27	0.64
1:D:354:ARG:HG2	1:D:354:ARG:NH1	2.10	0.63
1:D:373:GLN:HB2	4:D:621:HOH:O	1.99	0.63
1:D:84:ASP:OD1	3:D:502:A1JN3:N37	2.32	0.62
1:A:206:VAL:O	1:A:206:VAL:HG22	1.99	0.62
1:D:94:THR:HG23	1:D:253:TYR:CE2	2.33	0.62
1:A:42:PRO:HD2	4:A:623:HOH:O	1.99	0.61
1:A:50:ILE:CD1	1:A:59:VAL:HG21	2.30	0.61
1:D:118:ARG:NH2	1:D:118:ARG:HG2	2.04	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:64:ARG:NH2	1:D:317:VAL:HG21	2.15	0.60
1:D:3:GLY:N	1:D:4:GLU:CD	2.59	0.60
1:D:351:HIS:HB2	1:D:354:ARG:HG3	1.84	0.60
1:A:43:ALA:HB2	1:A:80:ASP:HB2	1.84	0.60
1:A:98:VAL:HG12	1:A:103:HIS:HB2	1.82	0.59
1:D:80:ASP:OD1	4:D:601:HOH:O	2.16	0.59
1:A:4:GLU:OE1	1:A:4:GLU:CA	2.50	0.59
1:D:77:ARG:HG2	1:D:77:ARG:HH11	1.67	0.59
1:D:85:ILE:HG23	1:D:85:ILE:O	2.03	0.58
1:D:312:PHE:CD1	1:D:312:PHE:N	2.69	0.58
1:A:403:ARG:NE	4:A:603:HOH:O	2.34	0.58
1:D:311:ARG:C	1:D:312:PHE:CD1	2.82	0.58
1:D:8:VAL:HG12	1:D:10:VAL:HG12	1.86	0.57
1:A:77:ARG:HG2	1:A:77:ARG:HH11	1.70	0.57
1:D:406:ARG:HH11	1:D:406:ARG:HG3	1.70	0.57
1:D:411:THR:HG23	1:D:413:ARG:NH2	2.20	0.56
1:A:98:VAL:HG13	1:A:102:ALA:HB3	1.88	0.56
1:D:64:ARG:NH1	1:D:317:VAL:HG21	2.20	0.56
1:D:67:LYS:O	1:D:68:ASP:C	2.44	0.56
1:A:389:ARG:C	1:A:391:PRO:HD3	2.30	0.56
1:A:390:PHE:N	1:A:391:PRO:HD3	2.21	0.56
1:A:70:ASP:HA	1:A:77:ARG:NH2	2.20	0.56
1:D:308:THR:HG23	1:D:332:VAL:HB	1.88	0.56
1:D:72:ALA:CA	1:D:312:PHE:HE2	2.18	0.56
1:A:125:ARG:O	1:A:129:ILE:HG13	2.06	0.55
1:A:70:ASP:HA	1:A:77:ARG:HH22	1.72	0.55
1:A:75:ALA:O	1:A:79:VAL:HG11	2.05	0.55
1:A:286:ALA:H	1:A:387:ARG:HH22	1.53	0.55
1:D:94:THR:HG23	1:D:253:TYR:HE2	1.72	0.55
1:D:70:ASP:HA	1:D:77:ARG:NH2	2.21	0.55
2:D:501:HEM:HBB2	2:D:501:HEM:CMB	2.35	0.55
1:D:354:ARG:HH11	1:D:354:ARG:CG	2.19	0.55
1:D:215:LEU:O	1:D:218:GLU:HB2	2.06	0.54
1:D:111:ALA:HB3	1:D:112:PRO:HD3	1.90	0.54
1:A:5:ALA:HB2	1:A:322:VAL:HG11	1.89	0.54
1:D:310:TRP:HB3	1:D:312:PHE:HE1	1.72	0.54
1:A:170:PRO:O	1:A:215:LEU:HD13	2.07	0.53
1:A:72:ALA:HB2	1:A:312:PHE:CZ	2.43	0.53
1:A:77:ARG:HH11	1:A:77:ARG:CG	2.20	0.53
1:A:25:GLU:O	1:A:26:VAL:C	2.49	0.53
1:D:265:THR:OG1	3:D:502:A1JN3:C34	2.57	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:THR:HA	1:A:322:VAL:O	2.09	0.53
1:D:72:ALA:CA	1:D:312:PHE:CE2	2.92	0.53
1:A:311:ARG:HH11	1:A:311:ARG:HG3	1.73	0.53
1:D:53:ASP:OD1	1:D:360:LEU:HD11	2.09	0.53
1:D:163:ILE:O	1:D:167:LEU:HD12	2.09	0.53
1:A:125:ARG:O	1:A:128:ALA:HB3	2.08	0.53
2:A:501:HEM:HBB2	2:A:501:HEM:HMB2	1.90	0.53
1:D:224:ARG:NE	1:D:248:ASP:OD1	2.41	0.53
1:D:311:ARG:C	1:D:312:PHE:HD1	2.17	0.53
1:D:411:THR:CG2	1:D:413:ARG:NH2	2.73	0.52
1:A:27:HIS:ND1	1:A:333:ASP:OD2	2.38	0.52
1:A:22:ARG:HG2	1:A:22:ARG:HH11	1.75	0.52
1:A:206:VAL:C	1:A:208:ASP:N	2.46	0.52
1:D:312:PHE:N	1:D:312:PHE:HD1	2.06	0.52
1:D:18:MET:O	1:D:403:ARG:HA	2.09	0.52
1:D:308:THR:CG2	1:D:332:VAL:HB	2.40	0.52
1:A:158:PHE:HB3	1:A:159:PRO:HD3	1.92	0.51
1:A:301:TYR:HE1	1:A:348:HIS:C	2.18	0.51
1:D:64:ARG:NH1	1:D:317:VAL:CG2	2.74	0.51
1:A:117:ARG:NH2	4:A:606:HOH:O	2.41	0.51
1:A:67:LYS:O	1:A:68:ASP:C	2.54	0.51
1:D:221:HIS:HB3	1:D:225:ARG:HH21	1.76	0.51
1:D:365:GLY:N	4:D:606:HOH:O	2.42	0.51
1:D:316:GLU:HG2	1:D:317:VAL:N	2.26	0.50
1:D:122:ARG:HH22	1:D:165:GLU:HG2	1.77	0.50
1:D:285:ASP:N	1:D:285:ASP:OD1	2.45	0.50
1:A:300:ARG:O	1:A:303:PRO:HD3	2.12	0.50
1:D:22:ARG:NH2	4:D:608:HOH:O	2.44	0.50
1:D:16:PRO:O	1:D:18:MET:HG2	2.12	0.49
1:D:94:THR:CG2	1:D:253:TYR:OH	2.60	0.49
1:A:90:LEU:HD12	4:A:669:HOH:O	2.11	0.49
1:D:297:GLU:CD	1:D:354:ARG:HD2	2.37	0.49
1:A:119:LEU:HD23	1:A:374:LEU:HD13	1.95	0.49
2:A:501:HEM:HBD1	2:A:501:HEM:HHA	1.95	0.49
1:D:286:ALA:HB1	1:D:290:ALA:HB3	1.95	0.49
1:A:320:GLY:N	4:A:604:HOH:O	2.39	0.48
1:D:159:PRO:HB2	1:D:263:HIS:HA	1.95	0.48
1:D:273:LEU:O	1:D:277:VAL:HG23	2.13	0.48
1:A:75:ALA:O	1:A:79:VAL:CG1	2.62	0.48
1:A:118:ARG:CZ	1:A:118:ARG:CB	2.92	0.48
1:A:411:THR:OG1	1:A:413:ARG:NH2	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:103:HIS:HE1	2:D:501:HEM:O2D	1.97	0.48
1:D:3:GLY:CA	1:D:4:GLU:CD	2.87	0.48
1:A:131:GLY:O	1:A:135:THR:OG1	2.28	0.47
1:A:215:LEU:O	1:A:218:GLU:HB3	2.14	0.47
1:D:307:TYR:CD1	1:D:333:ASP:HA	2.49	0.47
1:D:77:ARG:HH11	1:D:77:ARG:CG	2.27	0.47
1:D:26:VAL:O	1:D:29:ALA:HB3	2.13	0.47
1:A:21:GLN:HA	1:A:301:TYR:OH	2.15	0.47
1:A:108:ARG:HG3	1:A:108:ARG:HH11	1.80	0.47
1:D:8:VAL:HG12	1:D:10:VAL:CG1	2.44	0.47
1:D:286:ALA:H	1:D:387:ARG:NH2	2.12	0.47
1:D:389:ARG:NH2	4:D:610:HOH:O	2.47	0.47
1:A:111:ALA:HB3	1:A:112:PRO:HD3	1.97	0.47
1:A:206:VAL:O	1:A:206:VAL:HG23	2.13	0.46
1:A:159:PRO:HB2	1:A:263:HIS:HA	1.98	0.46
1:D:70:ASP:HA	1:D:77:ARG:HH22	1.81	0.46
1:A:351:HIS:O	1:A:354:ARG:HB2	2.16	0.46
1:A:409:ALA:O	3:A:502:A1JN3:O38	2.33	0.46
1:D:318:THR:HA	1:D:322:VAL:O	2.16	0.46
1:A:371:GLY:O	1:A:372:GLU:C	2.57	0.45
1:A:392:GLU:OE2	1:D:107:ARG:NE	2.49	0.45
1:D:393:ALA:HA	1:D:420:TRP:O	2.16	0.45
1:A:17:MET:SD	1:A:405:CYS:HB2	2.57	0.45
1:D:307:TYR:HD1	1:D:333:ASP:HA	1.81	0.45
1:D:158:PHE:HB3	1:D:159:PRO:CD	2.47	0.45
1:D:404:TRP:HB3	1:D:414:LEU:HD23	1.99	0.45
1:A:345:ASP:O	1:A:354:ARG:NH1	2.36	0.45
1:D:286:ALA:H	1:D:387:ARG:HH22	1.65	0.45
1:D:387:ARG:HH11	1:D:387:ARG:HD3	1.55	0.44
1:D:179:GLU:HG2	1:D:207:PRO:HB3	1.99	0.44
1:D:67:LYS:NZ	1:D:96:ILE:O	2.51	0.44
2:A:501:HEM:HBB2	2:A:501:HEM:CMB	2.47	0.44
1:D:94:THR:HG22	3:D:502:A1JN3:C28	2.48	0.44
1:D:10:VAL:O	1:D:39:VAL:CG2	2.66	0.43
1:D:318:THR:HG23	1:D:322:VAL:O	2.18	0.43
1:D:297:GLU:O	1:D:298:ALA:C	2.61	0.43
1:D:94:THR:O	1:D:95:LEU:C	2.62	0.43
1:A:98:VAL:CG1	1:A:102:ALA:HB3	2.47	0.43
1:D:333:ASP:O	1:D:337:THR:OG1	2.17	0.43
1:A:206:VAL:HA	1:A:207:PRO:HD3	1.73	0.43
1:D:25:GLU:O	1:D:26:VAL:C	2.62	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:77:ARG:HD2	1:D:77:ARG:N	2.33	0.43
1:D:211:ALA:O	1:D:215:LEU:CD2	2.54	0.43
3:A:502:A1JN3:O8	3:A:502:A1JN3:O36	2.37	0.43
1:A:107:ARG:NH2	1:A:368:TYR:O	2.34	0.42
1:A:22:ARG:HG2	1:A:22:ARG:NH1	2.35	0.42
1:D:371:GLY:O	1:D:372:GLU:C	2.63	0.42
1:D:50:ILE:HG23	1:D:55:LEU:HD23	2.02	0.42
1:D:207:PRO:O	1:D:208:ASP:C	2.63	0.42
1:A:184:LEU:HD21	1:A:209:THR:HB	2.01	0.42
1:A:389:ARG:C	1:A:390:PHE:CD1	2.97	0.42
1:D:67:LYS:O	1:D:69:PRO:N	2.53	0.42
1:D:318:THR:OG1	1:D:323:ARG:HG2	2.20	0.41
1:D:127:ALA:O	1:D:128:ALA:C	2.63	0.41
1:A:74:ALA:HA	1:A:77:ARG:HD3	2.02	0.41
1:A:317:VAL:HG12	1:A:318:THR:N	2.35	0.41
1:D:96:ILE:HG23	3:D:502:A1JN3:O19	2.21	0.41
1:A:12:PRO:HG3	1:A:39:VAL:HG21	2.02	0.41
1:A:156:TYR:O	1:A:159:PRO:HD2	2.20	0.41
2:A:501:HEM:HMD3	3:A:502:A1JN3:O19	2.20	0.41
1:A:143:ARG:NE	4:A:611:HOH:O	2.46	0.41
1:D:91:ARG:N	1:D:92:PRO:CD	2.84	0.41
1:D:300:ARG:O	1:D:303:PRO:HD3	2.21	0.41
1:A:335:GLU:O	1:A:339:THR:OG1	2.37	0.41
1:D:302:HIS:N	1:D:303:PRO:CD	2.84	0.41
1:D:376:GLN:O	1:D:377:LEU:C	2.64	0.41
1:D:60:LEU:HD23	1:D:311:ARG:NE	2.35	0.40
1:D:94:THR:HG21	1:D:253:TYR:OH	2.21	0.40
1:D:310:TRP:HB3	1:D:312:PHE:CE1	2.54	0.40
1:A:183:VAL:O	1:A:187:LEU:HG	2.21	0.40
1:A:387:ARG:HH11	1:A:387:ARG:HD2	1.58	0.40
1:D:36:VAL:HG21	1:D:319:ILE:HG22	2.03	0.40
1:D:104:ARG:O	1:D:105:ARG:C	2.63	0.40
1:D:317:VAL:HG22	1:D:318:THR:N	2.36	0.40
1:D:397:VAL:HG13	1:D:418:PRO:HG2	2.02	0.40
1:A:326:ARG:HH11	1:A:326:ARG:HD3	1.69	0.40
1:A:168:GLY:O	1:A:170:PRO:HD3	2.21	0.40
1:A:221:HIS:O	1:A:225:ARG:HG2	2.21	0.40
1:D:10:VAL:O	1:D:39:VAL:HG23	2.20	0.40
1:D:94:THR:HG23	1:D:253:TYR:CZ	2.56	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:225:ARG:NE	1:D:323:ARG:NH2[7_555]	2.12	0.08

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	403/457 (88%)	372 (92%)	27 (7%)	4 (1%)	12	17
1	D	403/457 (88%)	374 (93%)	26 (6%)	3 (1%)	18	25
All	All	806/914 (88%)	746 (93%)	53 (7%)	7 (1%)	14	20

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	81	ASP
1	D	207	PRO
1	A	79	VAL
1	A	207	PRO
1	A	341	GLY
1	A	356	SER
1	D	79	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	323/358 (90%)	304 (94%)	19 (6%)	18	27

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	322/358 (90%)	298 (92%)	24 (8%)	12	17
All	All	645/716 (90%)	602 (93%)	43 (7%)	15	21

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	17	MET
1	A	19	THR
1	A	77	ARG
1	A	81	ASP
1	A	98	VAL
1	A	118	ARG
1	A	193	GLN
1	A	206	VAL
1	A	227	ASP
1	A	230	THR
1	A	245	SER
1	A	265	THR
1	A	285	ASP
1	A	287	ASP
1	A	319	ILE
1	A	358	ARG
1	A	386	LEU
1	A	410	GLN
1	D	19	THR
1	D	67	LYS
1	D	77	ARG
1	D	79	VAL
1	D	81	ASP
1	D	94	THR
1	D	98	VAL
1	D	99	ASP
1	D	118	ARG
1	D	125	ARG
1	D	171	VAL
1	D	210	SER
1	D	214	SER
1	D	227	ASP
1	D	230	THR
1	D	245	SER
1	D	265	THR

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Mol	Chain	Res	Type
1	D	285	ASP
1	D	287	ASP
1	D	312	PHE
1	D	315	THR
1	D	345	ASP
1	D	354	ARG
1	D	411	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	250	GLN
1	D	103	HIS
1	D	344	HIS
1	D	367	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	A1JN3	A	502	2	40,42,42	0.98	2 (5%)	55,66,66	0.98	4 (7%)
3	A1JN3	D	502	2	40,42,42	0.67	1 (2%)	55,66,66	0.86	2 (3%)
2	HEM	D	501	1,3	50,50,50	1.99	15 (30%)	67,82,82	1.46	10 (14%)
2	HEM	A	501	1,3	50,50,50	2.28	16 (32%)	67,82,82	1.71	14 (20%)

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
3	A1JN3	A	502	2	-	4/12/58/58	0/5/5/5
3	A1JN3	D	502	2	-	0/12/58/58	0/5/5/5
2	HEM	D	501	1,3	-	4/14/54/54	-
2	HEM	A	501	1,3	-	3/14/54/54	-

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	FE-NC	8.46	2.23	1.95
2	A	501	HEM	FE-NA	5.28	2.12	1.95
2	D	501	HEM	C3C-C4C	-5.15	1.37	1.46
2	D	501	HEM	FE-ND	4.68	2.09	1.94
2	D	501	HEM	C1C-C2C	-4.22	1.37	1.45
2	A	501	HEM	C1C-C2C	-3.85	1.37	1.45
2	A	501	HEM	FE-NB	-3.74	1.83	1.94
2	D	501	HEM	C3B-C4B	-3.72	1.37	1.44
2	A	501	HEM	C4D-C3D	-3.71	1.38	1.45
2	A	501	HEM	C3C-C4C	-3.67	1.39	1.46
2	D	501	HEM	C4A-NA	-3.37	1.33	1.39
2	D	501	HEM	C1B-NB	-3.25	1.34	1.40
2	D	501	HEM	C4D-C3D	-3.13	1.39	1.45
2	D	501	HEM	C1C-NC	-3.13	1.33	1.39
2	D	501	HEM	C1D-C2D	-3.12	1.38	1.44
2	A	501	HEM	C4A-NA	-3.09	1.33	1.39
2	A	501	HEM	C1B-NB	-3.02	1.35	1.40
2	A	501	HEM	C3B-C4B	-3.01	1.39	1.44
2	A	501	HEM	C1A-C2A	-2.94	1.38	1.44
2	A	501	HEM	FE-ND	2.88	2.03	1.94
2	D	501	HEM	C1A-C2A	-2.83	1.38	1.44
2	A	501	HEM	C1D-C2D	-2.81	1.38	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	C1C-NC	-2.64	1.34	1.39
2	D	501	HEM	C1B-C2B	-2.52	1.39	1.44
2	D	501	HEM	C1A-NA	-2.46	1.35	1.39
2	D	501	HEM	C4A-C3A	-2.46	1.37	1.43
3	A	502	A1JN3	C10-C11	-2.44	1.50	1.53
2	D	501	HEM	C1D-ND	-2.42	1.34	1.38
3	A	502	A1JN3	O32-C15	2.30	1.42	1.36
3	D	502	A1JN3	C10-C11	-2.26	1.50	1.53
2	A	501	HEM	C3C-C2C	2.23	1.41	1.37
2	D	501	HEM	CAC-C3C	-2.22	1.41	1.47
2	A	501	HEM	C4C-NC	-2.19	1.35	1.39
2	A	501	HEM	O1D-CGD	2.08	1.28	1.22

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	HEM	CHD-C4C-NC	4.98	129.88	124.45
2	A	501	HEM	CHC-C4B-NB	4.93	129.73	124.42
2	A	501	HEM	CHC-C1C-NC	-4.42	119.64	124.45
2	A	501	HEM	C3B-C4B-NB	4.36	112.60	109.47
2	A	501	HEM	CHC-C4B-C3B	-3.72	117.66	125.07
2	A	501	HEM	C3C-C2C-C1C	-3.57	103.66	107.05
2	A	501	HEM	CHD-C4C-C3C	3.29	130.76	125.21
3	A	502	A1JN3	O7-C2-C1	-3.22	99.59	106.74
2	A	501	HEM	CHB-C4A-NA	-3.17	118.12	123.86
2	D	501	HEM	O1A-CGA-CBA	-3.13	113.17	123.09
2	A	501	HEM	CMB-C2B-C1B	3.00	129.72	125.03
2	A	501	HEM	CHD-C4C-NC	-2.99	121.20	124.45
2	A	501	HEM	C4B-C3B-C2B	-2.98	104.54	107.28
2	D	501	HEM	CAA-C2A-C1A	2.82	130.46	124.94
2	D	501	HEM	CAD-C3D-C4D	2.70	129.41	124.70
2	D	501	HEM	O2A-CGA-CBA	2.67	122.44	114.00
2	A	501	HEM	CMB-C2B-C3B	-2.56	122.24	128.43
3	A	502	A1JN3	O36-C11-C10	-2.55	103.06	108.40
3	D	502	A1JN3	C11-C10-C9	-2.46	109.62	114.58
2	D	501	HEM	CHD-C1D-C2D	2.43	128.87	125.03
2	D	501	HEM	C1D-C2D-C3D	2.40	109.50	106.98
2	D	501	HEM	CHD-C1D-ND	-2.39	121.85	124.42
3	A	502	A1JN3	O35-C33-C34	2.32	114.23	108.92
2	A	501	HEM	C4C-CHD-C1D	2.28	130.85	126.02
2	D	501	HEM	CHD-C4C-C3C	-2.24	121.44	125.21
3	D	502	A1JN3	O36-C11-C12	2.14	115.22	107.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	A1JN3	C11-C10-C9	-2.13	110.29	114.58
2	A	501	HEM	C3B-C2B-C1B	2.09	107.98	106.41
2	A	501	HEM	CMA-C3A-C4A	2.08	128.58	125.42
2	D	501	HEM	C4A-C3A-C2A	2.02	109.13	106.82

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	A1JN3	C10-C11-C33-C34
3	A	502	A1JN3	C12-C11-C33-C34
3	A	502	A1JN3	C12-C11-C33-O35
3	A	502	A1JN3	O36-C11-C33-C34
2	D	501	HEM	CAA-CBA-CGA-O2A
2	D	501	HEM	CAA-CBA-CGA-O1A
2	A	501	HEM	C2D-C3D-CAD-CBD
2	A	501	HEM	CAD-CBD-CGD-O1D
2	D	501	HEM	CAD-CBD-CGD-O2D
2	A	501	HEM	CAD-CBD-CGD-O2D
2	D	501	HEM	CAD-CBD-CGD-O1D

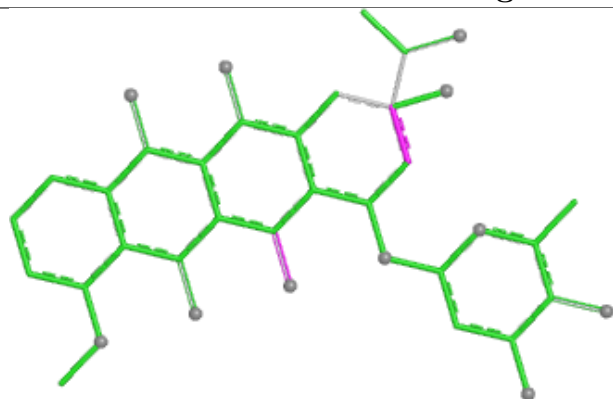
There are no ring outliers.

4 monomers are involved in 17 short contacts:

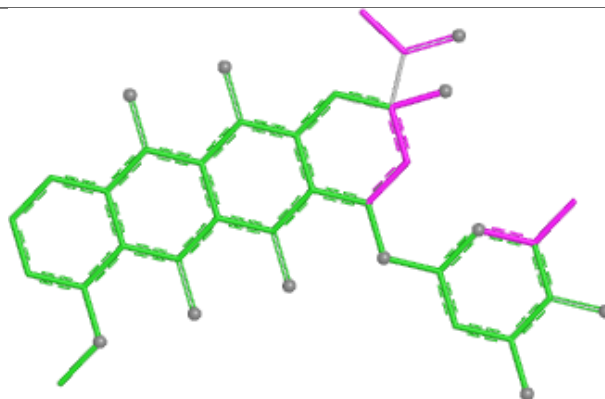
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	A1JN3	3	0
3	D	502	A1JN3	4	0
2	D	501	HEM	7	0
2	A	501	HEM	4	0

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.

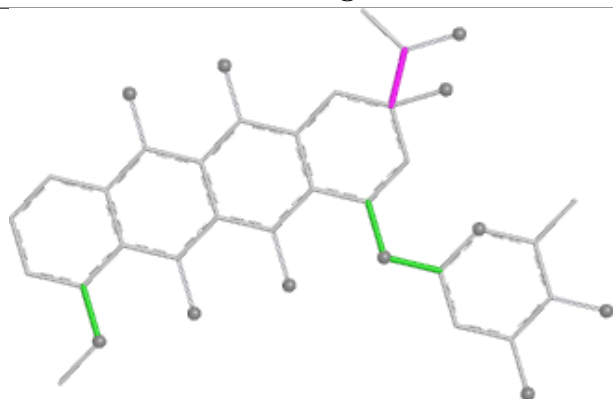
Ligand A1JN3 A 502



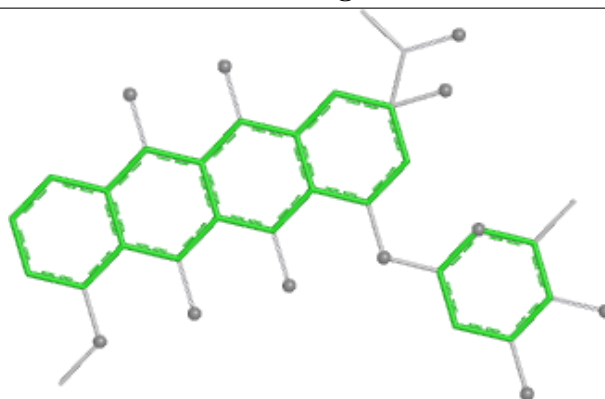
Bond lengths



Bond angles

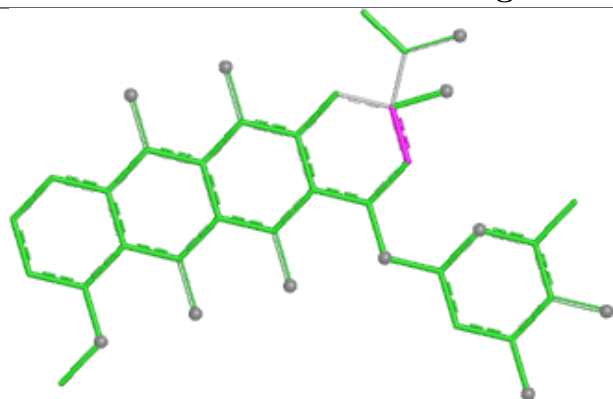


Torsions

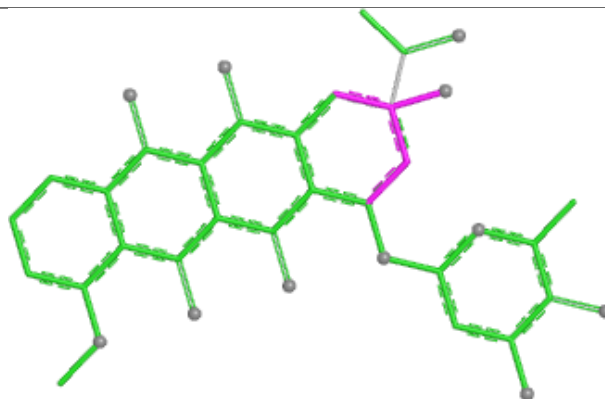


Rings

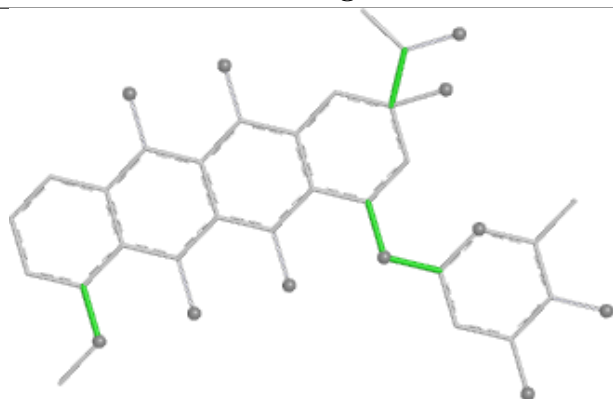
Ligand A1JN3 D 502



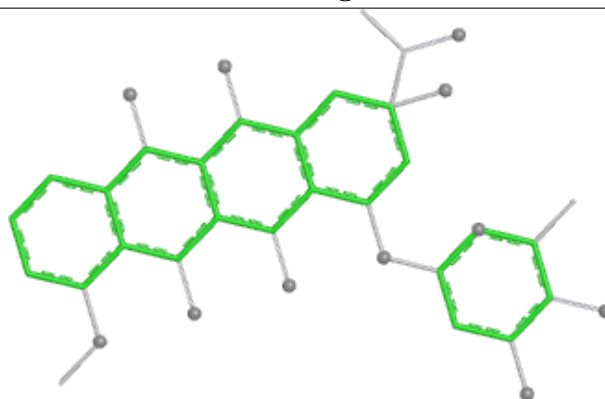
Bond lengths



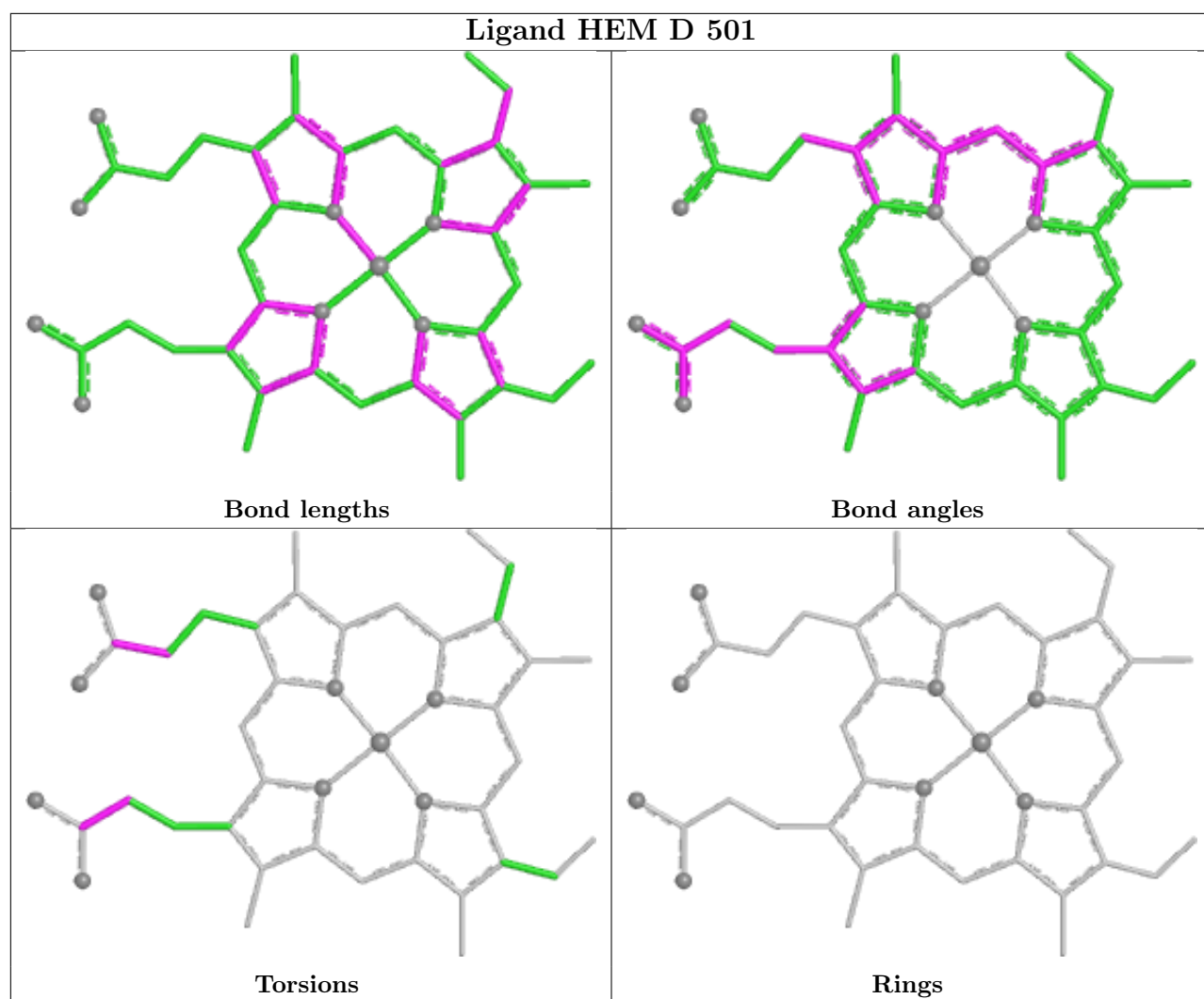
Bond angles

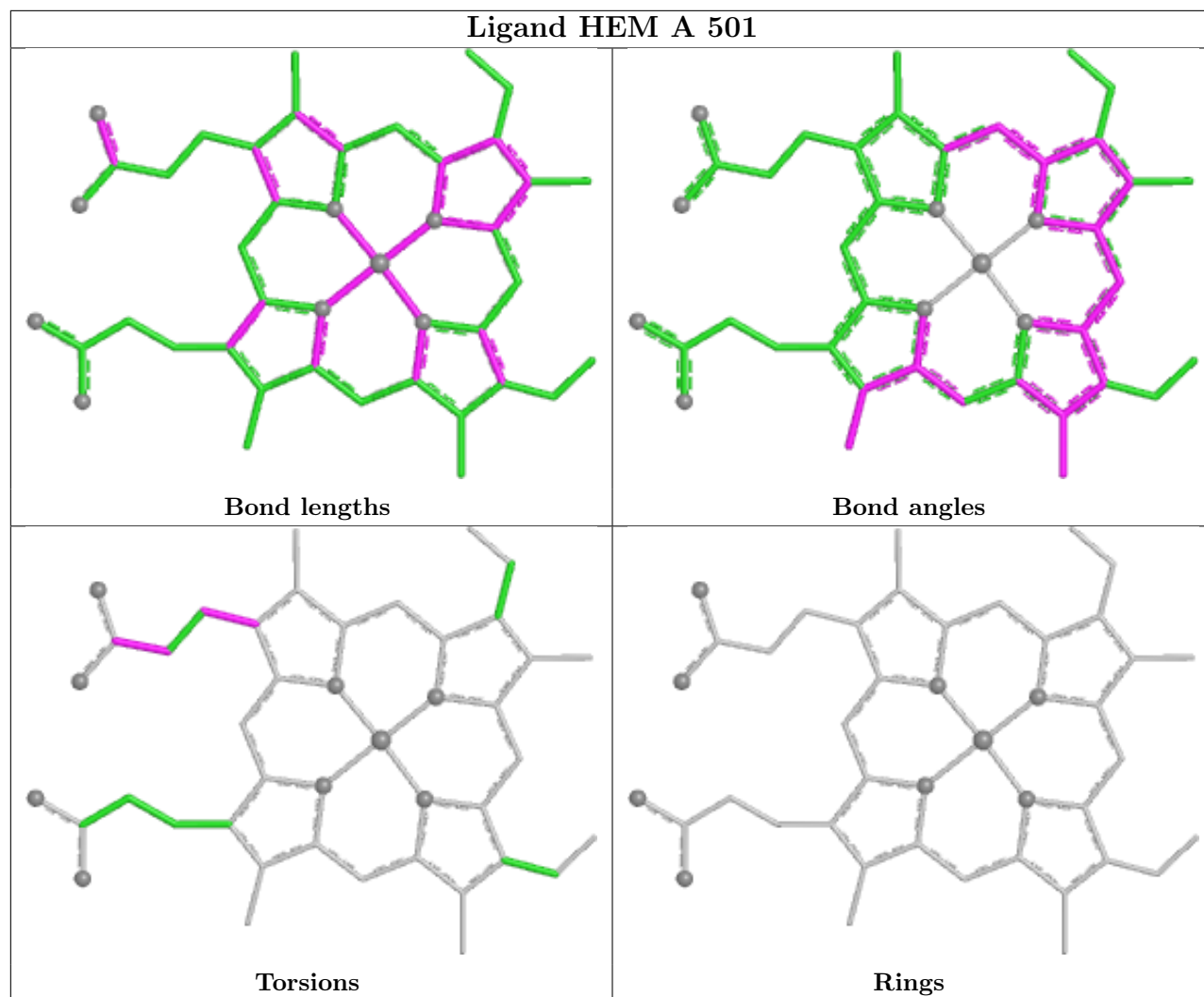


Torsions



Rings





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

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	407/457 (89%)	0.24	11 (2%) 56 57	39, 57, 87, 171	0
1	D	407/457 (89%)	0.86	56 (13%) 6 5	47, 76, 127, 186	0
All	All	814/914 (89%)	0.55	67 (8%) 17 17	39, 64, 114, 186	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	207	PRO	6.0
1	A	206	VAL	5.8
1	D	324	LEU	4.7
1	D	79	VAL	4.1
1	D	206	VAL	4.1
1	D	319	ILE	4.0
1	A	357	TRP	3.9
1	D	29	ALA	3.8
1	D	13	PHE	3.7
1	D	307	TYR	3.7
1	D	51	THR	3.5
1	D	322	VAL	3.3
1	D	35	PRO	3.3
1	D	333	ASP	3.2
1	D	32	GLU	3.1
1	D	59	VAL	3.1
1	D	192	PRO	3.0
1	D	317	VAL	3.0
1	D	61	ALA	3.0
1	D	75	ALA	3.0
1	D	65	PHE	3.0
1	D	50	ILE	2.8
1	D	11	ASP	2.8
1	D	355	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	230	THR	2.7
1	D	357	TRP	2.7
1	D	60	LEU	2.7
1	D	30	PHE	2.7
1	D	64	ARG	2.6
1	D	55	LEU	2.6
1	D	5	ALA	2.6
1	D	321	GLY	2.5
1	A	229	PRO	2.5
1	D	14	ALA	2.5
1	D	46	PRO	2.4
1	D	329	PRO	2.4
1	D	71	LEU	2.4
1	D	183	VAL	2.4
1	D	315	THR	2.3
1	D	26	VAL	2.3
1	D	313	ALA	2.3
1	D	207	PRO	2.3
1	D	230	THR	2.3
1	A	359	ARG	2.3
1	D	326	ARG	2.3
1	D	341	GLY	2.3
1	A	208	ASP	2.2
1	A	355	PRO	2.2
1	D	229	PRO	2.2
1	D	27	HIS	2.2
1	D	8	VAL	2.2
1	D	330	VAL	2.2
1	D	56	ALA	2.2
1	A	228	THR	2.2
1	A	368	TYR	2.2
1	D	301	TYR	2.2
1	D	320	GLY	2.1
1	D	187	LEU	2.1
1	D	36	VAL	2.1
1	D	3	GLY	2.1
1	D	325	PRO	2.1
1	D	6	PRO	2.1
1	D	243	PHE	2.0
1	D	85	ILE	2.0
1	D	57	ARG	2.0
1	A	319	ILE	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	73	PRO	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

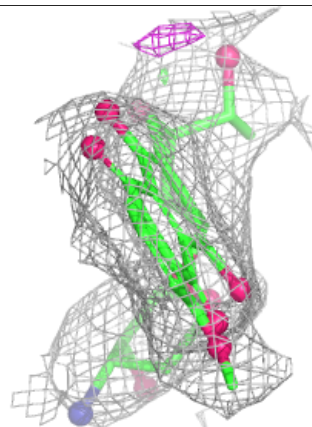
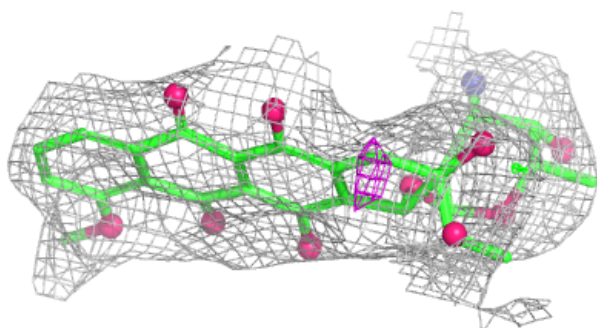
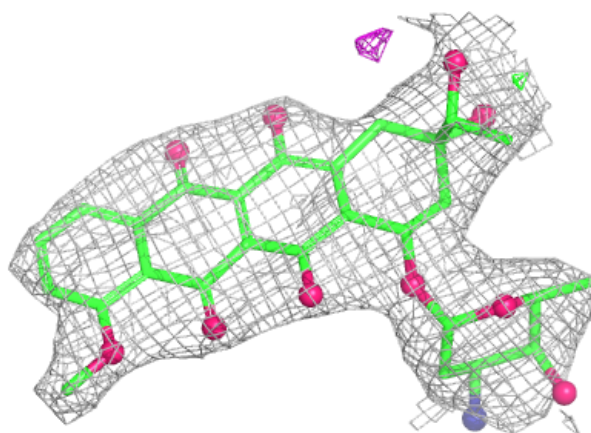
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	A1JN3	D	502	38/38	0.90	0.12	67,73,84,97	0
3	A1JN3	A	502	38/38	0.95	0.07	36,38,44,46	0
2	HEM	A	501	43/43	0.96	0.09	44,52,56,58	0
2	HEM	D	501	43/43	0.97	0.09	42,48,60,63	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.

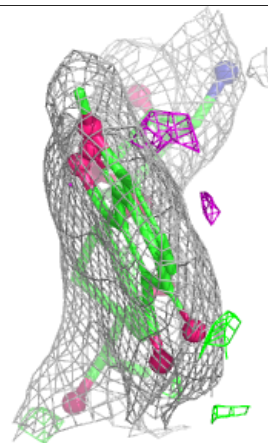
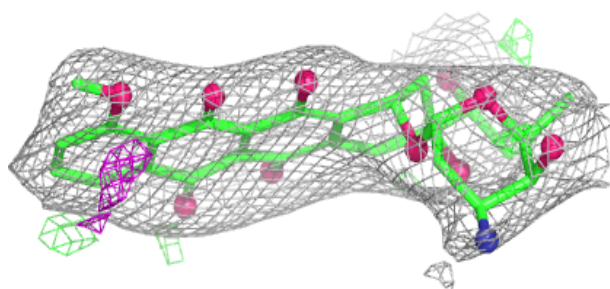
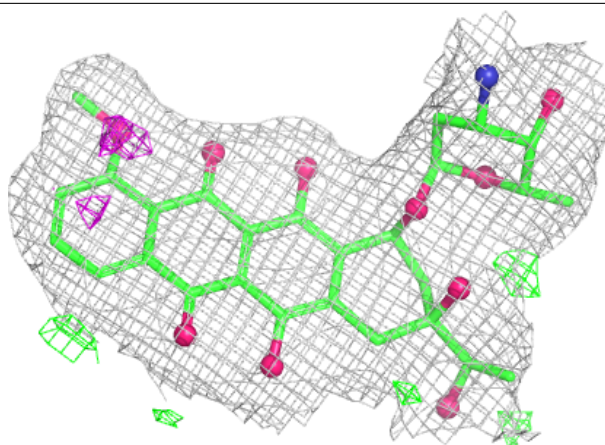
Electron density around A1JN3 D 502:

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



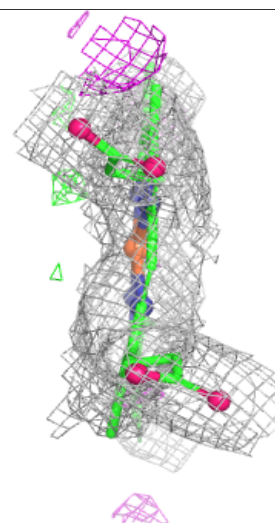
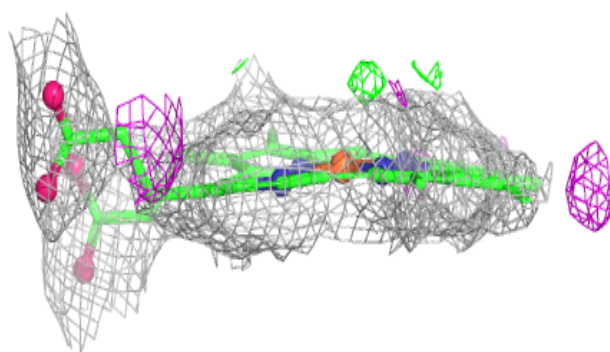
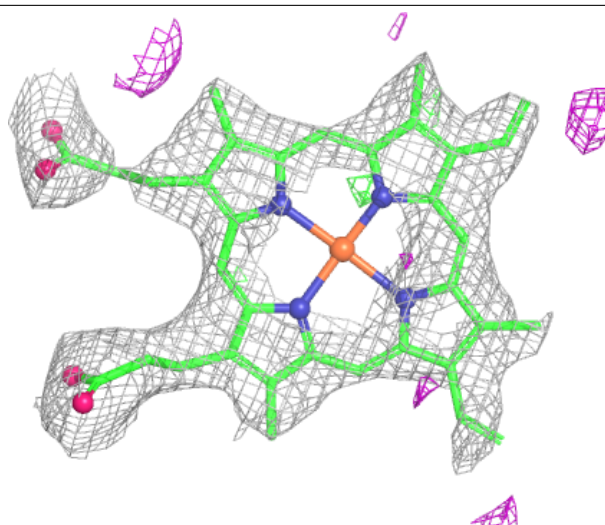
Electron density around A1JN3 A 502:

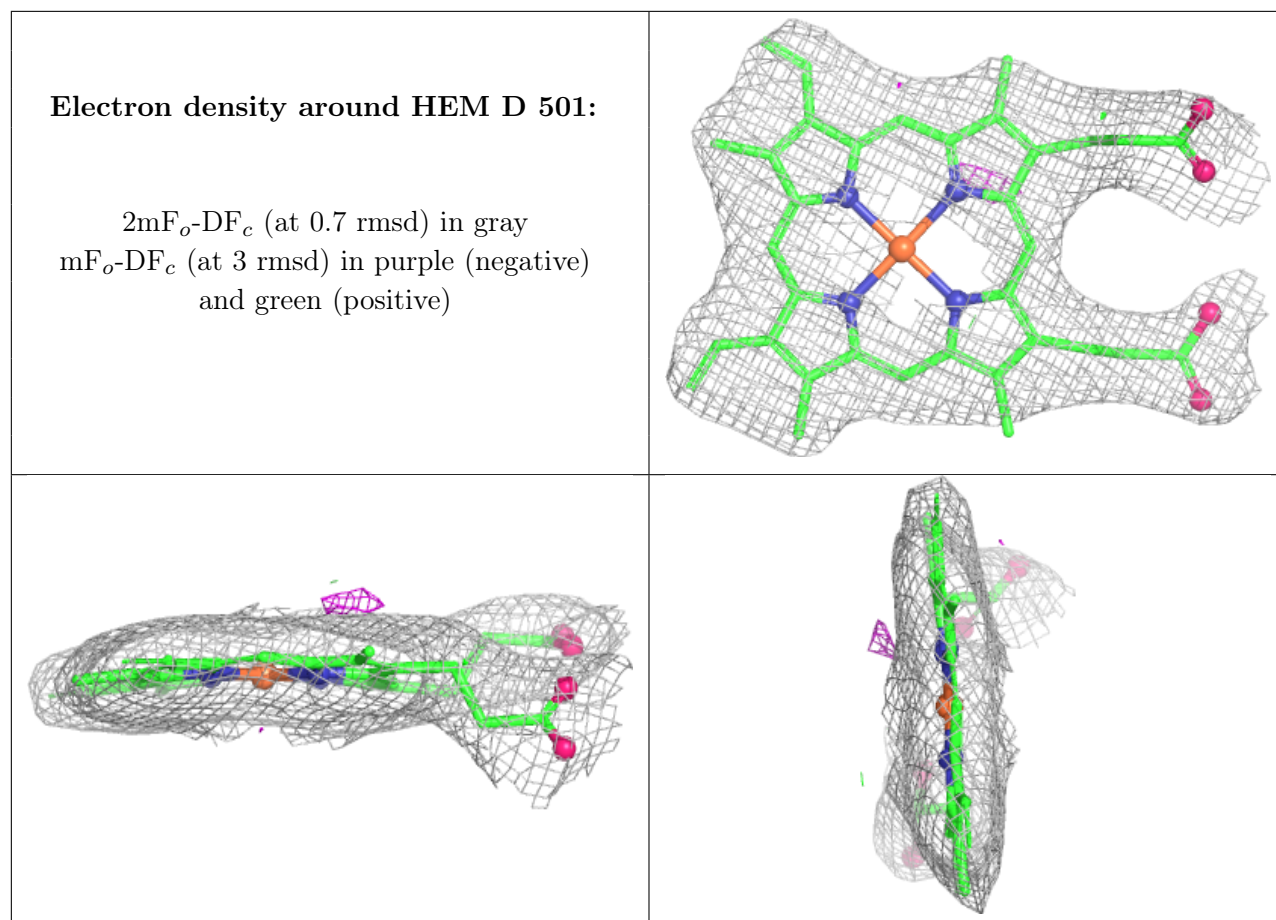
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 501:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.