



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:54 AM UTC

PDB ID : 4R20 / pdb_00004r20
Title : Zebra fish cytochrome P450 17A2 with Abiraterone
Authors : Pallan, P.S.; Egli, M.
Deposited on : 2014-08-08
Resolution : 2.86 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

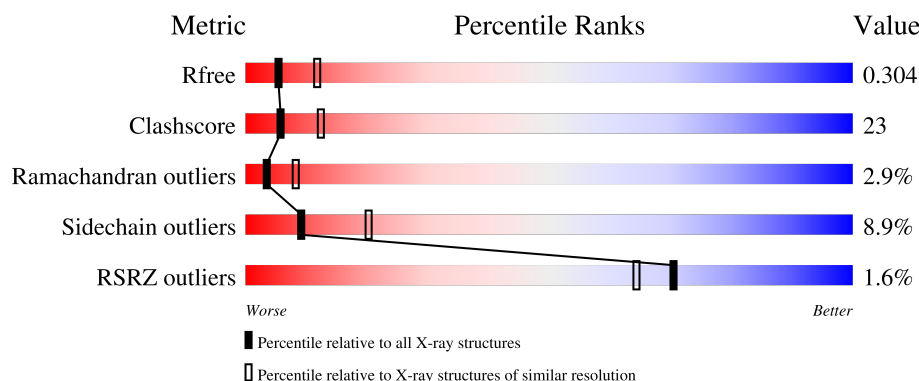
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.86 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	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	486	2% 57% 22% • • 16%
1	B	486	% 57% 22% 5% • 15%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6555 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 P450 family 17 polypeptide 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	406	Total	C	N	O	S	18	0	0
			3171	2011	563	582	15			
1	B	414	Total	C	N	O	S	5	0	0
			3227	2052	570	590	15			

There are 32 discrepancies between the modelled and reference sequences:

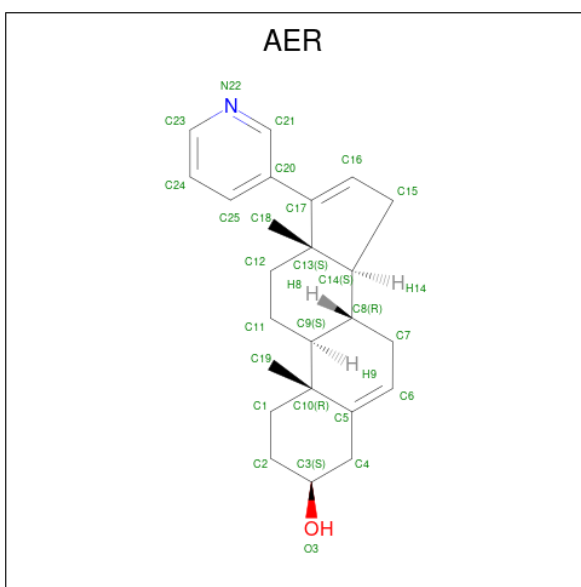
Chain	Residue	Modelled	Actual	Comment	Reference
A	16	MET	-	expression tag	UNP A7U483
A	17	ALA	-	expression tag	UNP A7U483
A	18	LYS	-	expression tag	UNP A7U483
A	19	LYS	-	expression tag	UNP A7U483
A	20	THR	-	expression tag	UNP A7U483
A	21	SER	-	expression tag	UNP A7U483
A	22	SER	-	expression tag	UNP A7U483
A	23	LYS	-	expression tag	UNP A7U483
A	24	GLY	-	expression tag	UNP A7U483
A	25	LYS	-	expression tag	UNP A7U483
A	496	HIS	-	expression tag	UNP A7U483
A	497	HIS	-	expression tag	UNP A7U483
A	498	HIS	-	expression tag	UNP A7U483
A	499	HIS	-	expression tag	UNP A7U483
A	500	HIS	-	expression tag	UNP A7U483
A	501	HIS	-	expression tag	UNP A7U483
B	16	MET	-	expression tag	UNP A7U483
B	17	ALA	-	expression tag	UNP A7U483
B	18	LYS	-	expression tag	UNP A7U483
B	19	LYS	-	expression tag	UNP A7U483
B	20	THR	-	expression tag	UNP A7U483
B	21	SER	-	expression tag	UNP A7U483
B	22	SER	-	expression tag	UNP A7U483
B	23	LYS	-	expression tag	UNP A7U483
B	24	GLY	-	expression tag	UNP A7U483

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Chain	Residue	Modelled	Actual	Comment	Reference
B	25	LYS	-	expression tag	UNP A7U483
B	496	HIS	-	expression tag	UNP A7U483
B	497	HIS	-	expression tag	UNP A7U483
B	498	HIS	-	expression tag	UNP A7U483
B	499	HIS	-	expression tag	UNP A7U483
B	500	HIS	-	expression tag	UNP A7U483
B	501	HIS	-	expression tag	UNP A7U483

- # HEM

- Molecule 3 is Abiraterone (CCD ID: AER) (formula: $C_{24}H_{31}NO$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			26	24	1	1		
3	B	1	Total	C	N	O	0	0
			26	24	1	1		

- Molecule 4 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Hg	0	0
			3	3		
4	B	2	Total	Hg	0	0
			2	2		

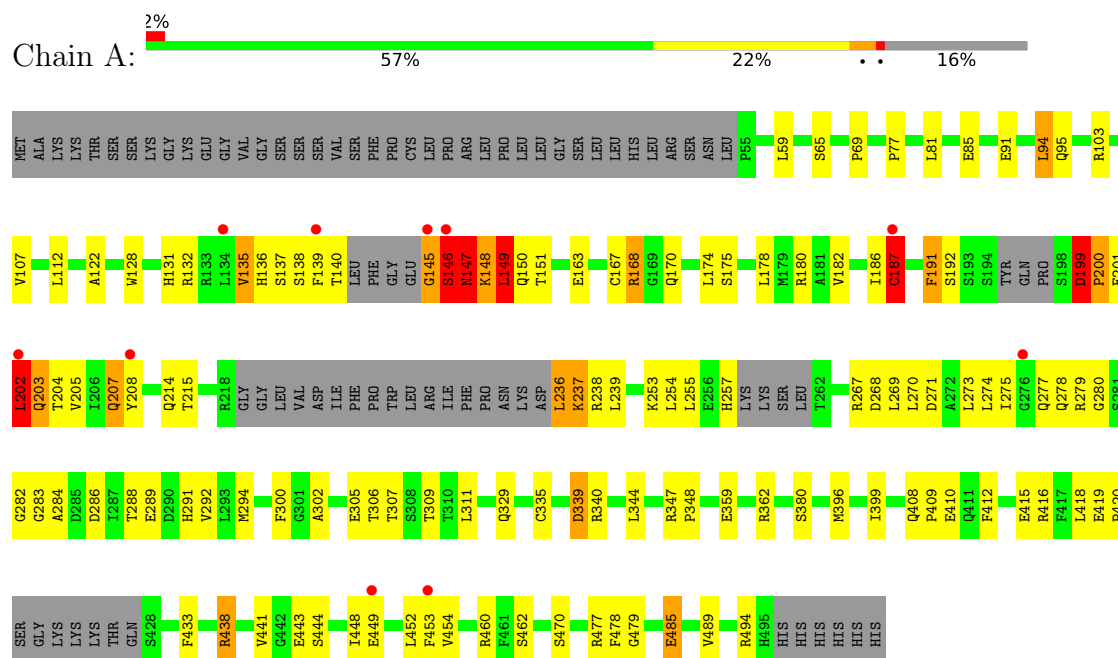
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	O	0	0
			3	3		
5	B	11	Total	O	0	0
			11	11		

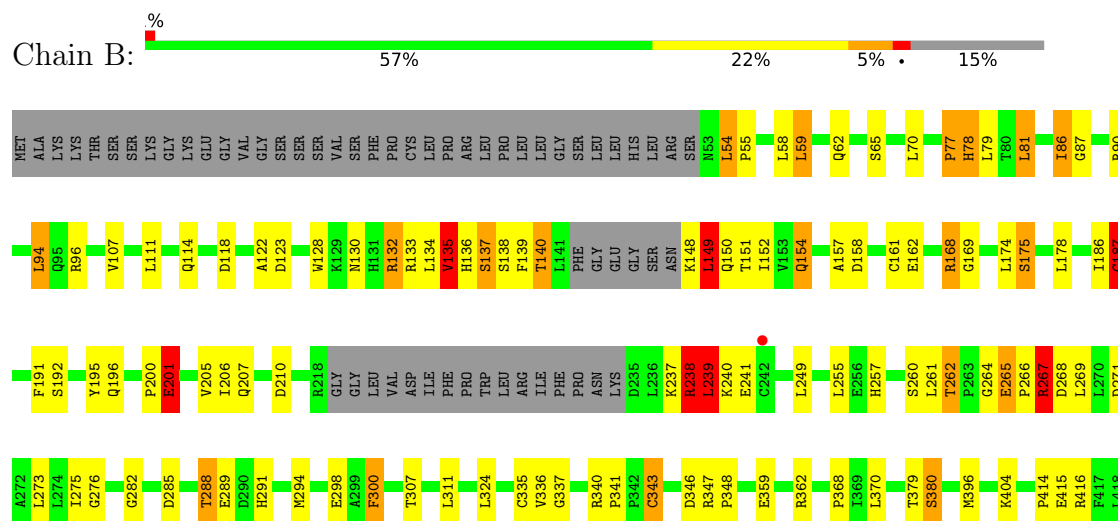
3 Residue-property plots

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: Cytochrome P450 family 17 polypeptide 2



- Molecule 1: Cytochrome P450 family 17 polypeptide 2



GLU	PRO	SER	GLY	LYS	LYS	LYS	THR	Q427	S428	S429	A435	R438	V439	C440	V441	E449	L452	F453	R456	S462	F463	S464	C465	P466	S467	E468	A469	S470	L471	F478	Q483	T490	V491	T492	P493	R494	H495	H496	HIS	HIS	HIS	HIS
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.01Å 79.77Å 95.24Å 90.00° 92.73° 90.00°	Depositor
Resolution (Å)	42.32 – 2.86 42.32 – 2.86	Depositor EDS
% Data completeness (in resolution range)	99.3 (42.32-2.86) 99.3 (42.32-2.86)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.98 (at 2.86Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.223 , 0.305 0.223 , 0.304	Depositor DCC
R_{free} test set	1005 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	79.5	Xtriage
Anisotropy	0.394	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 60.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.039 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6555	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.24% 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: HEM, HG, AER

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	0.69	1/3239 (0.0%)	1.05	8/4398 (0.2%)
1	B	0.76	1/3299 (0.0%)	1.08	17/4487 (0.4%)
All	All	0.73	2/6538 (0.0%)	1.06	25/8885 (0.3%)

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	7
1	B	0	3
All	All	0	10

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	464	SER	N-CA	5.08	1.52	1.45
1	A	187	CYS	CA-CB	5.04	1.61	1.53

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	267	ARG	CD-NE-CZ	11.11	139.95	124.40
1	B	267	ARG	NE-CZ-NH1	-10.56	110.94	121.50
1	A	187	CYS	N-CA-C	-9.86	100.53	111.28
1	A	146	SER	N-CA-C	-8.84	100.24	111.90
1	B	132	ARG	N-CA-C	-8.72	99.38	110.19
1	B	135	VAL	N-CA-C	-7.44	99.28	109.55
1	B	464	SER	N-CA-C	7.39	120.68	109.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	420	PRO	N-CA-CB	7.28	111.01	103.00
1	B	137	SER	N-CA-C	-6.68	103.92	111.07
1	B	187	CYS	N-CA-C	-6.66	104.05	111.71
1	B	300	PHE	N-CA-C	6.47	118.92	111.02
1	A	147	ASN	N-CA-C	-6.42	104.36	111.36
1	B	465	CYS	CA-C-N	-5.81	113.80	119.78
1	B	465	CYS	C-N-CA	-5.81	113.80	119.78
1	B	149	LEU	N-CA-C	5.64	119.03	111.24
1	A	199	ASP	C-N-CD	-5.45	108.61	120.60
1	B	465	CYS	CA-CB-SG	5.45	126.93	114.40
1	B	86	ILE	N-CA-C	5.32	116.69	110.62
1	B	107	VAL	CB-CA-C	-5.32	105.24	111.94
1	B	300	PHE	CB-CA-C	-5.32	102.47	110.92
1	A	202	LEU	CA-C-N	5.28	128.75	120.82
1	A	202	LEU	C-N-CA	5.28	128.75	120.82
1	A	454	VAL	N-CA-C	-5.24	107.66	111.90
1	B	238	ARG	CA-C-N	5.17	131.01	121.70
1	B	238	ARG	C-N-CA	5.17	131.01	121.70

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	145	GLY	Peptide
1	A	147	ASN	Peptide
1	A	149	LEU	Peptide
1	A	202	LEU	Peptide
1	A	236	LEU	Peptide
1	A	438	ARG	Sidechain
1	A	470	SER	Peptide
1	B	238	ARG	Peptide
1	B	267	ARG	Sidechain
1	B	428	SER	Peptide

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	3171	0	3171	138	0
1	B	3227	0	3221	151	0
2	A	43	0	30	6	0
2	B	43	0	30	7	0
3	A	26	0	31	4	0
3	B	26	0	31	4	0
4	A	3	0	0	0	0
4	B	2	0	0	1	0
5	A	3	0	0	0	0
5	B	11	0	0	0	0
All	All	6555	0	6514	301	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 23.

All (301) 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:A:205:VAL:HA	1:A:208:TYR:CE2	1.64	1.33
1:A:135:VAL:HG12	1:A:139:PHE:CE2	1.85	1.12
1:B:149:LEU:HB3	1:B:152:ILE:HG13	1.21	1.11
1:A:187:CYS:SG	1:A:300:PHE:HZ	1.77	1.06
1:B:465:CYS:SG	1:B:469:ALA:HB3	1.94	1.06
1:B:335:CYS:SG	4:B:604:HG:HG	1.86	0.94
1:B:264:GLY:C	1:B:265:GLU:OE1	2.11	0.93
1:A:146:SER:C	1:A:148:LYS:HB2	1.96	0.91
1:B:149:LEU:HB3	1:B:152:ILE:CG1	2.00	0.90
1:A:199:ASP:HB3	1:A:200:PRO:HA	1.54	0.90
1:B:139:PHE:O	1:B:140:THR:OG1	1.90	0.89
1:B:157:ALA:O	1:B:161:CYS:SG	2.32	0.88
1:B:192:SER:HB2	1:B:268:ASP:OD1	1.73	0.88
1:B:133:ARG:O	1:B:137:SER:HB3	1.75	0.86
1:B:149:LEU:N	1:B:150:GLN:HB3	1.91	0.84
1:A:205:VAL:HA	1:A:208:TYR:CD2	2.12	0.84
1:B:466:PRO:CD	1:B:469:ALA:HB2	2.08	0.83
1:B:148:LYS:C	1:B:150:GLN:HB3	2.03	0.83
1:A:103:ARG:HG2	1:A:438:ARG:NH2	1.94	0.83
1:A:103:ARG:HG2	1:A:438:ARG:HH21	1.43	0.82
1:A:267:ARG:HG2	1:A:271:ASP:OD2	1.80	0.82
1:A:199:ASP:HB3	1:A:200:PRO:CA	2.08	0.81
1:A:294:MET:HE2	1:A:294:MET:HA	1.62	0.81
1:A:410:GLU:N	1:A:410:GLU:OE1	2.13	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:LYS:O	1:B:149:LEU:HD22	1.80	0.81
1:B:466:PRO:N	1:B:469:ALA:HB2	1.96	0.81
1:A:146:SER:CA	1:A:148:LYS:HB2	2.11	0.81
1:B:467:SER:OG	1:B:468:GLU:OE2	1.99	0.81
1:B:466:PRO:HD2	1:B:469:ALA:HB2	1.62	0.80
1:B:136:HIS:HA	1:B:139:PHE:HB2	1.63	0.80
1:B:148:LYS:C	1:B:149:LEU:HD22	2.09	0.78
1:A:135:VAL:HG12	1:A:139:PHE:CZ	2.19	0.76
1:B:368:PRO:HD2	1:B:478:PHE:CE1	2.21	0.76
1:B:168:ARG:CD	1:B:169:GLY:N	2.49	0.75
1:B:168:ARG:CD	1:B:169:GLY:H	1.99	0.75
1:A:359:GLU:HG3	1:A:412:PHE:CE1	2.22	0.75
1:B:122:ALA:O	1:B:438:ARG:NH2	2.20	0.74
1:B:149:LEU:H	1:B:151:THR:H	1.35	0.74
1:A:149:LEU:HA	1:A:150:GLN:C	2.11	0.74
1:A:199:ASP:HB2	1:A:200:PRO:HB3	1.70	0.74
1:B:187:CYS:SG	1:B:300:PHE:HZ	2.11	0.73
1:A:205:VAL:CA	1:A:208:TYR:CE2	2.60	0.73
1:A:146:SER:HA	1:A:148:LYS:HB2	1.70	0.73
1:B:134:LEU:O	1:B:138:SER:N	2.23	0.72
1:B:149:LEU:N	1:B:149:LEU:HD13	2.05	0.71
1:A:277:GLN:NE2	1:A:289:GLU:HG3	2.06	0.71
1:A:135:VAL:CG1	1:A:139:PHE:CE2	2.70	0.71
1:B:168:ARG:HD3	1:B:169:GLY:H	1.54	0.71
1:B:134:LEU:HD22	1:B:273:LEU:HD23	1.73	0.71
1:A:274:LEU:O	1:A:278:GLN:HG3	1.91	0.70
1:A:146:SER:OG	1:A:149:LEU:CD1	2.41	0.69
1:B:149:LEU:HA	1:B:151:THR:N	2.08	0.69
1:A:187:CYS:HG	1:A:300:PHE:HZ	0.75	0.69
2:A:601:HEM:HHC	2:A:601:HEM:HBB2	1.73	0.69
1:A:191:PHE:HE2	1:A:300:PHE:CE1	2.11	0.68
1:B:466:PRO:HD2	1:B:469:ALA:CB	2.23	0.68
1:A:236:LEU:HD13	1:A:237:LYS:HB2	1.75	0.68
1:B:168:ARG:HD2	1:B:169:GLY:N	2.09	0.68
1:A:149:LEU:HA	1:A:151:THR:N	2.09	0.68
1:B:150:GLN:HG3	1:B:151:THR:N	2.09	0.68
1:B:132:ARG:NH2	1:B:439:VAL:O	2.27	0.67
1:B:149:LEU:N	1:B:151:THR:H	1.91	0.67
1:A:267:ARG:CG	1:A:271:ASP:OD2	2.43	0.67
1:B:139:PHE:C	1:B:140:THR:HG1	1.99	0.67
1:B:134:LEU:C	1:B:134:LEU:HD23	2.21	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:ARG:O	1:B:137:SER:CB	2.42	0.66
1:A:202:LEU:HA	1:A:204:THR:N	2.10	0.66
1:B:368:PRO:HD2	1:B:478:PHE:HE1	1.59	0.65
1:A:128:TRP:CE2	1:A:438:ARG:NH1	2.65	0.65
1:B:294:MET:HE2	1:B:294:MET:HA	1.79	0.65
1:B:135:VAL:O	1:B:136:HIS:HB3	1.94	0.65
1:A:128:TRP:NE1	1:A:438:ARG:NH1	2.44	0.65
1:A:199:ASP:CB	1:A:200:PRO:HB3	2.26	0.64
1:A:103:ARG:CG	1:A:438:ARG:NH2	2.61	0.64
2:B:601:HEM:HHD	2:B:601:HEM:HBC2	1.79	0.64
1:B:238:ARG:N	1:B:239:LEU:HB3	2.13	0.63
1:A:122:ALA:O	1:A:438:ARG:NH2	2.28	0.63
1:B:196:GLN:H	1:B:200:PRO:HG2	1.65	0.62
1:B:136:HIS:CA	1:B:139:PHE:HB2	2.29	0.62
1:B:311:LEU:HD11	1:B:453:PHE:CE1	2.34	0.62
1:A:478:PHE:CD1	1:A:478:PHE:C	2.78	0.62
1:A:146:SER:OG	1:A:149:LEU:HD11	1.99	0.61
1:A:203:GLN:O	1:A:207:GLN:N	2.24	0.61
1:A:236:LEU:HA	1:A:237:LYS:CB	2.31	0.61
1:B:239:LEU:HD13	1:B:239:LEU:C	2.26	0.61
1:A:478:PHE:C	1:A:478:PHE:HD1	2.08	0.61
1:A:146:SER:O	1:A:149:LEU:HD13	2.00	0.61
1:B:128:TRP:CH2	1:B:132:ARG:HD2	2.35	0.61
1:B:132:ARG:O	1:B:133:ARG:HB3	2.00	0.61
1:A:199:ASP:CB	1:A:200:PRO:CA	2.78	0.60
1:B:149:LEU:H	1:B:151:THR:HG23	1.65	0.60
1:B:149:LEU:HA	1:B:150:GLN:C	2.26	0.60
1:A:236:LEU:HA	1:A:237:LYS:HB2	1.84	0.60
1:B:134:LEU:O	1:B:134:LEU:HD23	2.01	0.60
1:A:91:GLU:HA	1:A:95:GLN:HB3	1.84	0.60
1:B:136:HIS:N	1:B:139:PHE:H	2.00	0.60
1:B:186:ILE:HG21	1:B:300:PHE:HA	1.83	0.60
1:A:289:GLU:HA	1:A:292:VAL:HG23	1.83	0.60
2:B:601:HEM:NB	3:B:602:AER:H23	2.16	0.60
1:B:134:LEU:CD2	1:B:273:LEU:HD23	2.32	0.60
1:A:200:PRO:HB2	1:A:202:LEU:O	2.02	0.59
1:A:339:ASP:O	1:A:340:ARG:C	2.46	0.59
1:A:191:PHE:HE2	1:A:300:PHE:CZ	2.20	0.59
1:B:136:HIS:O	1:B:139:PHE:O	2.20	0.59
1:A:362:ARG:NH2	1:A:409:PRO:O	2.35	0.59
1:A:191:PHE:CE2	1:A:300:PHE:CE1	2.91	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:PHE:HB3	1:A:253:LYS:HG2	1.85	0.58
1:A:329:GLN:HG2	1:A:494:ARG:NH2	2.18	0.58
1:B:264:GLY:O	1:B:265:GLU:OE1	2.19	0.58
1:A:311:LEU:HD11	1:A:453:PHE:HE1	1.68	0.58
1:A:415:GLU:HA	1:A:418:LEU:HG	1.84	0.58
2:A:601:HEM:HBC2	2:A:601:HEM:HHD	1.85	0.58
1:B:151:THR:HA	1:B:154:GLN:HG3	1.86	0.58
1:A:282:GLY:HA2	1:A:283:GLY:C	2.28	0.58
1:A:477:ARG:NH2	1:A:485:GLU:HG2	2.19	0.57
2:A:601:HEM:C1D	3:A:602:AER:H21	2.39	0.57
1:A:136:HIS:O	1:A:137:SER:C	2.46	0.57
1:A:275:ILE:O	1:A:279:ARG:HG3	2.05	0.57
1:A:204:THR:O	1:A:208:TYR:CD2	2.58	0.56
1:B:187:CYS:CB	1:B:195:TYR:HE2	2.19	0.56
1:B:265:GLU:OE1	1:B:265:GLU:N	2.38	0.56
2:B:601:HEM:C1D	3:B:602:AER:H21	2.39	0.56
1:B:267:ARG:NH1	1:B:271:ASP:OD2	2.33	0.56
1:B:266:PRO:O	1:B:267:ARG:HG2	2.06	0.56
1:A:187:CYS:SG	1:A:300:PHE:CZ	2.70	0.55
1:A:145:GLY:C	1:A:147:ASN:N	2.55	0.55
1:A:135:VAL:CG1	1:A:139:PHE:CZ	2.89	0.55
1:A:174:LEU:HD23	1:A:489:VAL:HG11	1.89	0.54
1:A:146:SER:HA	1:A:148:LYS:CB	2.37	0.54
1:B:288:THR:OG1	1:B:289:GLU:N	2.41	0.54
1:A:478:PHE:HD1	1:A:479:GLY:N	2.05	0.54
1:B:187:CYS:SG	1:B:300:PHE:CZ	2.95	0.54
1:A:135:VAL:O	1:A:139:PHE:CD2	2.61	0.54
1:B:130:ASN:ND2	1:B:285:ASP:HB3	2.23	0.54
1:B:266:PRO:O	1:B:271:ASP:OD2	2.25	0.53
1:B:261:LEU:HD11	1:B:275:ILE:CD1	2.38	0.53
1:A:131:HIS:O	1:A:135:VAL:HG23	2.08	0.53
1:A:307:THR:HG21	1:A:449:GLU:OE1	2.09	0.53
2:B:601:HEM:ND	3:B:602:AER:H21	2.24	0.53
2:A:601:HEM:NB	3:A:602:AER:H23	2.23	0.53
1:B:168:ARG:HD3	1:B:169:GLY:N	2.15	0.53
1:B:441:VAL:HG22	2:B:601:HEM:HMD2	1.90	0.53
2:B:601:HEM:C1B	3:B:602:AER:H23	2.44	0.53
1:A:137:SER:O	1:A:140:THR:HG23	2.09	0.53
1:A:168:ARG:O	1:A:170:GLN:HG2	2.09	0.53
1:A:174:LEU:CD2	1:A:489:VAL:HG11	2.38	0.53
1:A:215:THR:HG21	1:A:239:LEU:HB2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:LEU:HD11	1:B:81:LEU:HG	1.89	0.53
1:A:277:GLN:CD	1:A:289:GLU:HG3	2.34	0.52
1:A:302:ALA:O	1:A:306:THR:OG1	2.22	0.52
1:B:307:THR:HG21	1:B:449:GLU:OE1	2.09	0.52
2:A:601:HEM:ND	3:A:602:AER:H21	2.24	0.52
1:B:134:LEU:HD11	1:B:276:GLY:HA3	1.91	0.52
1:A:202:LEU:HA	1:A:203:GLN:C	2.34	0.52
1:A:200:PRO:CB	1:A:202:LEU:O	2.57	0.52
1:A:433:PHE:CD2	1:A:443:GLU:HG3	2.45	0.52
1:B:267:ARG:HG2	1:B:271:ASP:OD2	2.10	0.51
1:A:236:LEU:HD13	1:A:237:LYS:HD2	1.91	0.51
2:A:601:HEM:C1B	3:A:602:AER:H23	2.46	0.51
1:B:150:GLN:NE2	1:B:154:GLN:OE1	2.43	0.51
1:B:261:LEU:HD11	1:B:275:ILE:HD11	1.93	0.51
1:B:55:PRO:HD2	1:B:58:LEU:HD12	1.93	0.51
1:A:329:GLN:CG	1:A:494:ARG:NH2	2.73	0.51
1:B:133:ARG:O	1:B:137:SER:N	2.39	0.51
1:A:359:GLU:CG	1:A:412:PHE:CE1	2.93	0.50
1:A:277:GLN:CD	1:A:289:GLU:CG	2.84	0.50
1:B:187:CYS:SG	1:B:195:TYR:HE2	2.35	0.50
1:B:158:ASP:OD1	1:B:456:ARG:NH1	2.44	0.50
1:B:267:ARG:HH11	1:B:271:ASP:CG	2.16	0.50
1:A:149:LEU:O	1:A:149:LEU:HD22	2.11	0.50
1:A:236:LEU:HA	1:A:237:LYS:CG	2.42	0.50
1:A:128:TRP:CH2	1:A:132:ARG:HD2	2.47	0.49
1:A:135:VAL:O	1:A:138:SER:HB3	2.12	0.49
1:A:148:LYS:O	1:A:151:THR:CB	2.60	0.49
1:A:199:ASP:CB	1:A:200:PRO:CB	2.89	0.49
1:A:236:LEU:CD1	1:A:237:LYS:HB2	2.41	0.49
1:A:329:GLN:HB3	1:A:494:ARG:HH22	1.76	0.49
1:A:203:GLN:N	1:A:203:GLN:OE1	2.46	0.49
1:A:289:GLU:O	1:A:292:VAL:N	2.42	0.49
1:B:187:CYS:HB3	1:B:195:TYR:HE2	1.77	0.49
1:B:195:TYR:CE1	1:B:201:GLU:OE2	2.66	0.49
1:B:111:LEU:HD13	1:B:239:LEU:HD12	1.95	0.49
1:B:118:ASP:OD2	1:B:298:GLU:OE2	2.30	0.49
1:B:492:THR:HB	1:B:493:PRO:CD	2.42	0.49
1:A:149:LEU:HD22	1:A:149:LEU:C	2.37	0.49
1:B:347:ARG:HB2	1:B:348:PRO:HD3	1.94	0.49
1:B:149:LEU:HB3	1:B:152:ILE:CD1	2.42	0.49
1:A:205:VAL:HG22	1:A:208:TYR:CE2	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:LYS:O	1:B:148:LYS:HG3	2.13	0.48
1:A:167:CYS:O	1:A:170:GLN:HB2	2.12	0.48
1:B:178:LEU:HD22	1:B:311:LEU:HD13	1.95	0.48
1:B:191:PHE:HE2	1:B:300:PHE:HE1	1.61	0.48
1:A:149:LEU:HD13	1:A:149:LEU:H	1.78	0.48
1:A:174:LEU:HD13	1:A:178:LEU:HD11	1.95	0.48
1:B:78:HIS:ND1	1:B:78:HIS:N	2.61	0.48
1:B:135:VAL:HG23	1:B:139:PHE:HD2	1.79	0.48
1:B:136:HIS:H	1:B:139:PHE:H	1.62	0.48
1:B:191:PHE:CE2	1:B:300:PHE:HE1	2.31	0.48
1:A:202:LEU:O	1:A:202:LEU:HD12	2.14	0.48
1:A:145:GLY:HA2	1:A:147:ASN:H	1.78	0.47
1:B:288:THR:HG23	1:B:291:HIS:CE1	2.48	0.47
1:A:163:GLU:OE1	1:A:180:ARG:NH1	2.48	0.47
1:B:468:GLU:N	1:B:468:GLU:CD	2.73	0.47
1:A:199:ASP:HB3	1:A:200:PRO:CB	2.44	0.47
1:A:204:THR:HG23	1:A:205:VAL:N	2.29	0.47
1:B:86:ILE:O	1:B:90:ARG:HG3	2.14	0.47
1:B:148:LYS:C	1:B:149:LEU:HD13	2.39	0.47
1:B:148:LYS:HA	1:B:149:LEU:HD13	1.96	0.47
1:A:146:SER:OG	1:A:149:LEU:HD12	2.12	0.47
1:A:148:LYS:O	1:A:151:THR:N	2.47	0.47
1:A:282:GLY:HA2	1:A:283:GLY:O	2.15	0.46
1:A:200:PRO:HB2	1:A:202:LEU:C	2.40	0.46
1:A:236:LEU:HB3	1:A:239:LEU:H	1.80	0.46
1:B:134:LEU:C	1:B:134:LEU:CD2	2.88	0.46
1:B:195:TYR:HE1	1:B:201:GLU:OE2	1.97	0.46
1:A:168:ARG:C	1:A:168:ARG:CD	2.89	0.46
1:B:149:LEU:CA	1:B:151:THR:N	2.76	0.46
2:B:601:HEM:HHC	2:B:601:HEM:HBB2	1.97	0.46
1:A:254:LEU:HA	1:A:270:LEU:HD21	1.96	0.46
1:A:288:THR:OG1	1:A:289:GLU:N	2.49	0.46
1:B:154:GLN:NE2	1:B:456:ARG:HH21	2.14	0.46
1:A:347:ARG:HB2	1:A:348:PRO:HD3	1.98	0.46
1:B:238:ARG:CB	1:B:239:LEU:HB2	2.45	0.46
1:B:135:VAL:CG2	1:B:139:PHE:HD2	2.29	0.46
1:B:168:ARG:HD2	1:B:169:GLY:H	1.70	0.46
1:A:128:TRP:CZ2	1:A:132:ARG:HD2	2.51	0.46
1:B:249:LEU:HD23	1:B:249:LEU:O	2.15	0.46
1:B:86:ILE:HG23	1:B:87:GLY:N	2.31	0.46
1:A:294:MET:HE2	1:A:294:MET:CA	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:237:LYS:HA	1:B:238:ARG:CB	2.46	0.45
1:B:149:LEU:HA	1:B:152:ILE:H	1.82	0.45
1:B:257:HIS:HA	1:B:260:SER:HB2	1.98	0.45
1:B:158:ASP:O	1:B:162:GLU:N	2.37	0.45
1:B:238:ARG:O	1:B:241:GLU:HG3	2.17	0.45
1:A:269:LEU:O	1:A:273:LEU:HG	2.16	0.45
1:B:135:VAL:HG23	1:B:139:PHE:CD2	2.52	0.45
1:B:237:LYS:CA	1:B:238:ARG:CB	2.95	0.45
1:B:130:ASN:ND2	1:B:285:ASP:CB	2.78	0.45
1:A:128:TRP:NE1	1:A:438:ARG:HH12	2.14	0.45
1:B:174:LEU:O	1:B:175:SER:CB	2.64	0.45
1:A:433:PHE:CG	1:A:443:GLU:HG3	2.51	0.45
1:B:135:VAL:HA	1:B:138:SER:CB	2.47	0.45
1:A:277:GLN:NE2	1:A:289:GLU:CG	2.78	0.44
1:A:283:GLY:HA2	1:A:284:ALA:C	2.42	0.44
1:A:182:VAL:O	1:A:186:ILE:HG12	2.17	0.44
1:B:465:CYS:HB2	1:B:471:LEU:HD23	1.99	0.44
1:B:174:LEU:O	1:B:175:SER:HB2	2.18	0.44
1:B:261:LEU:HD12	1:B:262:THR:N	2.33	0.44
1:A:280:GLY:HA3	1:A:286:ASP:CG	2.43	0.44
1:A:282:GLY:HA3	1:A:286:ASP:H	1.83	0.44
1:B:174:LEU:C	1:B:174:LEU:HD12	2.42	0.44
1:B:370:LEU:HD11	1:B:396:MET:HE3	1.99	0.44
1:A:128:TRP:CZ2	1:A:438:ARG:NH1	2.81	0.44
1:B:111:LEU:CD2	1:B:240:LYS:HG2	2.48	0.43
1:B:379:THR:OG1	1:B:380:SER:N	2.49	0.43
1:A:204:THR:O	1:A:208:TYR:HD2	2.01	0.43
1:B:128:TRP:CZ3	1:B:132:ARG:HD2	2.52	0.43
1:B:77:PRO:HA	1:B:78:HIS:HA	1.65	0.43
1:B:54:LEU:HD23	1:B:59:LEU:HD23	2.00	0.43
1:A:136:HIS:O	1:A:140:THR:HG23	2.18	0.43
1:A:204:THR:CG2	1:A:205:VAL:N	2.81	0.43
1:A:257:HIS:NE2	1:A:268:ASP:OD2	2.37	0.43
1:B:467:SER:HG	1:B:468:GLU:CD	2.27	0.43
1:A:344:LEU:HD21	1:A:448:ILE:HG23	2.01	0.42
1:B:148:LYS:C	1:B:149:LEU:CD2	2.88	0.42
1:A:192:SER:HB2	1:A:268:ASP:OD1	2.19	0.42
1:B:187:CYS:SG	1:B:195:TYR:CE2	3.12	0.42
1:B:324:LEU:HD21	1:B:414:PRO:HG2	2.01	0.42
1:B:493:PRO:O	1:B:494:ARG:HG3	2.18	0.42
1:B:257:HIS:O	1:B:261:LEU:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:SER:O	1:B:267:ARG:NH2	2.52	0.42
1:B:466:PRO:O	1:B:469:ALA:CB	2.68	0.42
1:B:239:LEU:C	1:B:239:LEU:CD1	2.93	0.42
1:B:336:VAL:HG12	1:B:337:GLY:N	2.35	0.42
1:A:203:GLN:CD	1:A:203:GLN:H	2.27	0.42
1:A:305:GLU:O	1:A:309:THR:OG1	2.28	0.42
1:B:111:LEU:HD21	1:B:240:LYS:HG2	2.01	0.41
1:B:466:PRO:HG3	1:B:490:THR:OG1	2.20	0.41
1:A:69:PRO:HB2	1:A:85:GLU:HB2	2.01	0.41
1:B:136:HIS:HA	1:B:139:PHE:CB	2.43	0.41
1:B:340:ARG:HA	1:B:341:PRO:HD2	1.86	0.41
1:A:288:THR:CG2	1:A:291:HIS:CD2	3.04	0.41
1:A:359:GLU:OE1	1:A:359:GLU:HA	2.19	0.41
1:B:149:LEU:CA	1:B:150:GLN:C	2.94	0.41
1:B:239:LEU:HD13	1:B:240:LYS:N	2.36	0.41
1:B:340:ARG:NH2	1:B:346:ASP:OD2	2.54	0.41
1:B:341:PRO:HB3	1:B:456:ARG:NH1	2.35	0.41
1:B:94:LEU:HD21	1:B:435:ALA:HB1	2.03	0.41
1:B:205:VAL:HG11	1:B:300:PHE:CE2	2.56	0.41
1:A:205:VAL:HA	1:A:208:TYR:CZ	2.41	0.41
1:A:408:GLN:HA	1:A:410:GLU:OE2	2.20	0.41
1:A:148:LYS:O	1:A:151:THR:HB	2.21	0.40
1:A:396:MET:HA	1:A:399:ILE:HG12	2.02	0.40
1:B:168:ARG:HD3	1:B:168:ARG:HA	1.97	0.40
1:B:359:GLU:OE2	1:B:362:ARG:NH1	2.53	0.40
1:A:94:LEU:O	1:A:95:GLN:C	2.63	0.40
1:A:254:LEU:HD12	1:A:274:LEU:HD11	2.04	0.40
1:B:187:CYS:HB3	1:B:195:TYR:CE2	2.55	0.40
1:B:340:ARG:NH2	1:B:343:CYS:SG	2.91	0.40
1:A:201:GLU:CG	1:A:202:LEU:N	2.85	0.40
1:B:135:VAL:O	1:B:136:HIS:CB	2.65	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	394/486 (81%)	367 (93%)	19 (5%)	8 (2%)	6	13
1	B	406/486 (84%)	365 (90%)	26 (6%)	15 (4%)	2	5
All	All	800/972 (82%)	732 (92%)	45 (6%)	23 (3%)	3	8

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	148	LYS
1	A	199	ASP
1	A	200	PRO
1	A	419	GLU
1	B	79	LEU
1	B	175	SER
1	B	201	GLU
1	B	238	ARG
1	B	239	LEU
1	B	429	SER
1	B	492	THR
1	B	493	PRO
1	A	175	SER
1	B	428	SER
1	B	495	HIS
1	B	54	LEU
1	A	339	ASP
1	B	282	GLY
1	B	470	SER
1	B	77	PRO
1	B	140	THR
1	A	77	PRO
1	A	237	LYS

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	354/427 (83%)	327 (92%)	27 (8%)	12	26
1	B	357/427 (84%)	321 (90%)	36 (10%)	7	15
All	All	711/854 (83%)	648 (91%)	63 (9%)	9	20

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	LEU
1	A	65	SER
1	A	81	LEU
1	A	94	LEU
1	A	107	VAL
1	A	112	LEU
1	A	135	VAL
1	A	146	SER
1	A	149	LEU
1	A	168	ARG
1	A	187	CYS
1	A	191	PHE
1	A	202	LEU
1	A	203	GLN
1	A	207	GLN
1	A	214	GLN
1	A	238	ARG
1	A	255	LEU
1	A	335	CYS
1	A	380	SER
1	A	416	ARG
1	A	441	VAL
1	A	444	SER
1	A	452	LEU
1	A	460	ARG
1	A	462	SER
1	A	485	GLU
1	B	59	LEU
1	B	62	GLN
1	B	65	SER
1	B	78	HIS
1	B	81	LEU
1	B	94	LEU
1	B	96	ARG
1	B	114	GLN

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Mol	Chain	Res	Type
1	B	123	ASP
1	B	135	VAL
1	B	149	LEU
1	B	154	GLN
1	B	168	ARG
1	B	187	CYS
1	B	201	GLU
1	B	206	ILE
1	B	207	GLN
1	B	210	ASP
1	B	239	LEU
1	B	255	LEU
1	B	262	THR
1	B	265	GLU
1	B	267	ARG
1	B	269	LEU
1	B	288	THR
1	B	343	CYS
1	B	380	SER
1	B	404	LYS
1	B	415	GLU
1	B	416	ARG
1	B	452	LEU
1	B	462	SER
1	B	465	CYS
1	B	470	SER
1	B	483	GLN
1	B	495	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	114	GLN
1	A	248	GLN
1	A	277	GLN
1	A	291	HIS
1	A	495	HIS
1	B	130	ASN
1	B	150	GLN
1	B	154	GLN
1	B	203	GLN
1	B	207	GLN

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Mol	Chain	Res	Type
1	B	395	ASN
1	B	405	HIS

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 ⓘ

Of 9 ligands modelled in this entry, 5 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	AER	A	602	2	30,30,30	2.34	2 (6%)	47,47,47	2.63	18 (38%)
3	AER	B	602	2	30,30,30	2.41	2 (6%)	47,47,47	2.65	16 (34%)
2	HEM	B	601	1,3	50,50,50	1.68	9 (18%)	67,82,82	2.32	31 (46%)
2	HEM	A	601	1,3	50,50,50	1.65	12 (24%)	67,82,82	1.92	16 (23%)

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	AER	A	602	2	-	2/4/62/62	0/5/5/5
3	AER	B	602	2	-	1/4/62/62	0/5/5/5
2	HEM	B	601	1,3	-	7/14/54/54	-
2	HEM	A	601	1,3	-	2/14/54/54	-

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	602	AER	C16-C17	12.16	1.54	1.33
3	A	602	AER	C16-C17	11.98	1.54	1.33
2	B	601	HEM	FE-NB	4.76	2.09	1.94
2	A	601	HEM	FE-NB	4.58	2.09	1.94
2	B	601	HEM	C4D-ND	-4.37	1.32	1.40
2	A	601	HEM	C4D-ND	-4.08	1.33	1.40
2	B	601	HEM	FE-NC	3.63	2.07	1.95
2	B	601	HEM	C1B-NB	-3.47	1.34	1.40
2	A	601	HEM	FE-NC	3.33	2.06	1.95
2	A	601	HEM	C1B-NB	-3.16	1.34	1.40
2	A	601	HEM	FE-ND	-3.01	1.85	1.94
2	A	601	HEM	C1C-C2C	-2.87	1.39	1.45
2	B	601	HEM	C1D-ND	-2.72	1.33	1.38
3	B	602	AER	C13-C17	2.66	1.56	1.53
2	A	601	HEM	C1D-ND	-2.64	1.33	1.38
2	A	601	HEM	C1A-NA	-2.40	1.35	1.39
2	B	601	HEM	C3D-C2D	-2.35	1.31	1.36
2	A	601	HEM	C3B-C4B	2.29	1.49	1.44
3	A	602	AER	C13-C17	2.21	1.55	1.53
2	B	601	HEM	C3B-C4B	2.20	1.49	1.44
2	B	601	HEM	C3C-C4C	-2.14	1.42	1.46
2	A	601	HEM	C4B-NB	-2.14	1.34	1.38
2	A	601	HEM	C4C-NC	-2.13	1.35	1.39
2	B	601	HEM	CHA-C4D	-2.06	1.34	1.38
2	A	601	HEM	CHA-C1A	-2.03	1.34	1.39

All (81) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	602	AER	C15-C16-C17	-10.03	103.86	112.56
3	B	602	AER	C20-C17-C16	-9.01	113.46	125.08
3	A	602	AER	C20-C17-C16	-8.75	113.79	125.08
3	A	602	AER	C15-C16-C17	-7.77	105.82	112.56
2	B	601	HEM	C3B-C4B-NB	-5.42	105.58	109.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	HEM	C1B-NB-C4B	4.76	110.84	105.21
2	A	601	HEM	CHA-C4D-ND	4.63	130.09	124.37
3	A	602	AER	C18-C13-C12	-4.47	105.99	111.10
2	A	601	HEM	CHD-C1D-ND	4.38	129.14	124.42
2	A	601	HEM	C1B-NB-C4B	4.37	110.38	105.21
3	A	602	AER	C15-C14-C8	-4.28	117.09	121.63
2	B	601	HEM	CHB-C1B-NB	4.27	129.65	124.37
3	A	602	AER	C13-C17-C20	-4.23	120.91	125.39
2	B	601	HEM	CHA-C4D-C3D	-4.22	117.44	125.23
3	A	602	AER	C15-C14-C13	4.22	107.00	104.05
2	B	601	HEM	CAA-C2A-C3A	4.10	136.27	127.07
2	B	601	HEM	CAA-C2A-C1A	-4.10	116.94	124.94
2	B	601	HEM	CHD-C1D-ND	3.95	128.68	124.42
3	B	602	AER	C13-C17-C16	-3.85	106.24	109.66
2	A	601	HEM	CHC-C4B-NB	3.84	128.56	124.42
3	A	602	AER	C13-C17-C16	-3.82	106.26	109.66
2	A	601	HEM	CHA-C4D-C3D	-3.78	118.27	125.23
3	B	602	AER	C15-C14-C8	-3.72	117.69	121.63
2	B	601	HEM	CBA-CAA-C2A	3.65	122.64	112.53
2	A	601	HEM	CHB-C1B-NB	3.64	128.87	124.37
2	B	601	HEM	CAA-CBA-CGA	-3.53	104.30	113.67
2	B	601	HEM	C3D-C4D-ND	3.52	114.03	110.17
3	A	602	AER	C14-C13-C17	3.44	102.42	99.50
2	A	601	HEM	C3B-C4B-NB	-3.41	107.02	109.47
2	B	601	HEM	CHA-C4D-ND	3.36	128.53	124.37
3	A	602	AER	C12-C11-C9	3.36	118.84	113.14
3	B	602	AER	C14-C8-C9	-3.24	104.85	109.09
2	A	601	HEM	CHD-C1D-C2D	-3.21	119.96	125.03
2	B	601	HEM	CMD-C2D-C1D	3.20	130.04	125.03
3	B	602	AER	C14-C13-C17	3.19	102.20	99.50
3	B	602	AER	C12-C13-C14	-3.17	103.97	108.88
3	A	602	AER	C3-C4-C5	-3.17	107.01	112.05
2	B	601	HEM	CAD-C3D-C4D	3.06	130.03	124.70
2	B	601	HEM	CAB-C3B-C2B	-3.02	118.60	128.43
2	B	601	HEM	C3B-C2B-C1B	3.00	108.66	106.41
2	A	601	HEM	CAA-CBA-CGA	-2.98	105.77	113.67
2	B	601	HEM	O2A-CGA-O1A	-2.93	115.80	123.33
3	B	602	AER	C13-C17-C20	-2.89	122.33	125.39
2	A	601	HEM	CHC-C1C-NC	2.88	127.59	124.45
2	B	601	HEM	CHD-C1D-C2D	-2.83	120.56	125.03
2	A	601	HEM	CBA-CAA-C2A	2.82	120.33	112.53
3	B	602	AER	C12-C13-C17	2.80	122.53	119.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	HEM	CAC-C3C-C4C	2.73	131.33	124.82
2	B	601	HEM	O2D-CGD-O1D	-2.72	116.34	123.33
3	A	602	AER	C12-C13-C17	2.70	122.43	119.30
3	B	602	AER	C18-C13-C12	-2.68	108.04	111.10
2	A	601	HEM	O2A-CGA-CBA	2.67	122.45	114.00
2	A	601	HEM	CAC-C3C-C4C	2.65	131.16	124.82
2	B	601	HEM	C4C-NC-C1C	2.58	110.03	105.82
3	B	602	AER	C24-C25-C20	-2.56	117.85	120.36
3	A	602	AER	C2-C3-C4	-2.54	106.72	110.29
2	B	601	HEM	CHC-C4B-NB	2.48	127.09	124.42
2	B	601	HEM	O2A-CGA-CBA	2.46	121.77	114.00
3	B	602	AER	C3-C4-C5	-2.44	108.16	112.05
2	B	601	HEM	O2D-CGD-CBD	2.43	121.69	114.00
2	B	601	HEM	CAB-C3B-C4B	2.40	134.99	124.39
3	B	602	AER	C12-C11-C9	2.38	117.19	113.14
3	B	602	AER	C15-C14-C13	2.38	105.71	104.05
3	B	602	AER	C4-C5-C10	2.32	119.39	116.42
3	B	602	AER	C13-C14-C8	2.29	115.55	113.13
3	A	602	AER	C12-C13-C14	-2.29	105.33	108.88
3	A	602	AER	C19-C10-C9	2.28	114.22	111.66
2	A	601	HEM	O2D-CGD-CBD	2.27	121.18	114.00
2	A	601	HEM	C3B-C2B-C1B	2.26	108.11	106.41
3	A	602	AER	C4-C5-C10	2.19	119.23	116.42
2	B	601	HEM	CMC-C2C-C3C	-2.16	123.19	128.43
2	B	601	HEM	CHC-C1C-NC	2.15	126.80	124.45
2	B	601	HEM	C4D-C3D-C2D	-2.13	103.79	106.89
2	B	601	HEM	CAC-C3C-C2C	-2.13	121.52	128.43
2	B	601	HEM	C1A-CHA-C4D	-2.11	121.28	126.25
3	A	602	AER	C10-C9-C8	-2.10	109.64	112.71
2	B	601	HEM	CHA-C1A-NA	2.09	127.65	123.86
3	A	602	AER	O3-C3-C2	-2.06	104.78	110.17
2	B	601	HEM	CMB-C2B-C3B	-2.06	123.45	128.43
3	A	602	AER	O3-C3-C4	2.05	114.33	109.71
2	A	601	HEM	C4A-CHB-C1B	-2.01	121.51	126.25

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	601	HEM	C3A-C2A-CAA-CBA
2	B	601	HEM	C1A-C2A-CAA-CBA
2	A	601	HEM	C4B-C3B-CAB-CBB

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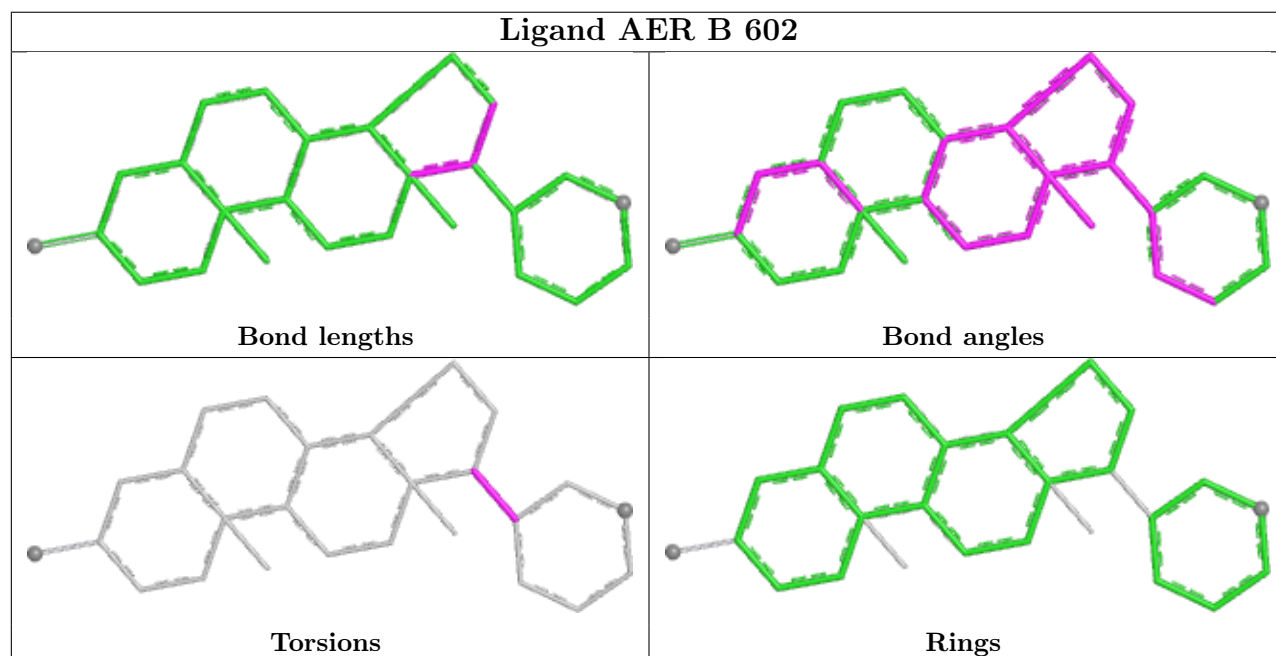
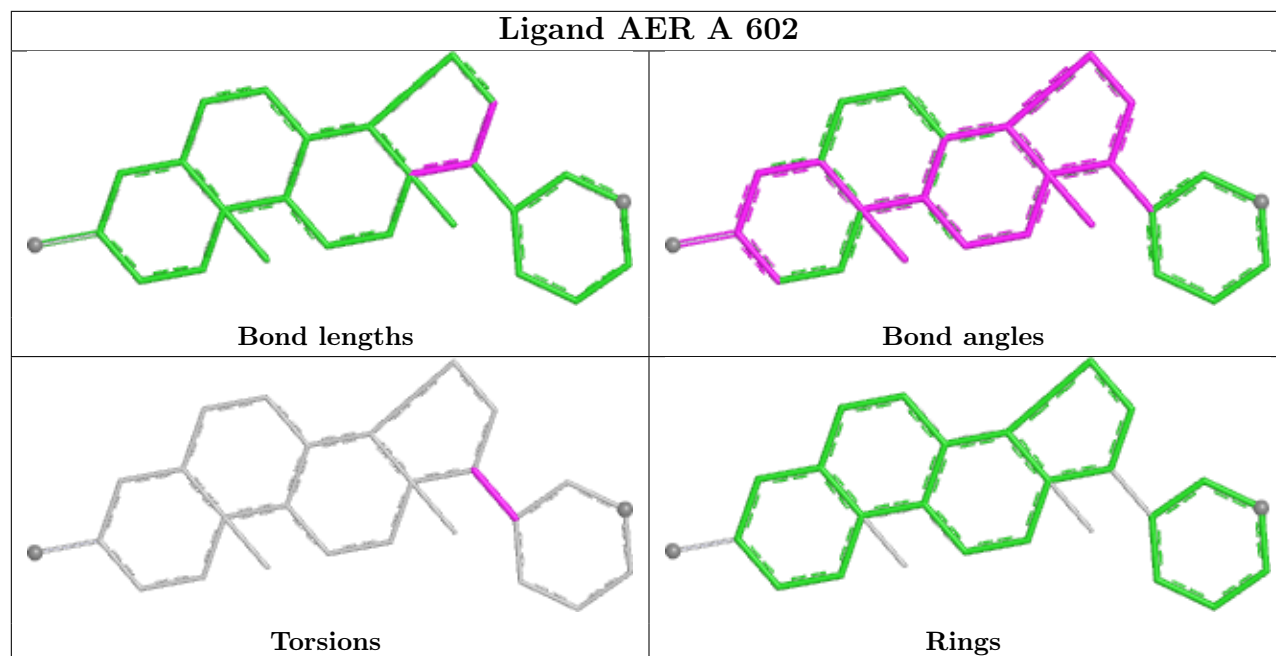
Mol	Chain	Res	Type	Atoms
2	A	601	HEM	C4C-C3C-CAC-CBC
2	B	601	HEM	C4B-C3B-CAB-CBB
2	B	601	HEM	C4C-C3C-CAC-CBC
2	B	601	HEM	CAA-CBA-CGA-O1A
2	B	601	HEM	CAA-CBA-CGA-O2A
3	A	602	AER	C13-C17-C20-C25
3	B	602	AER	C13-C17-C20-C25
3	A	602	AER	C13-C17-C20-C21
2	B	601	HEM	CAD-CBD-CGD-O2D

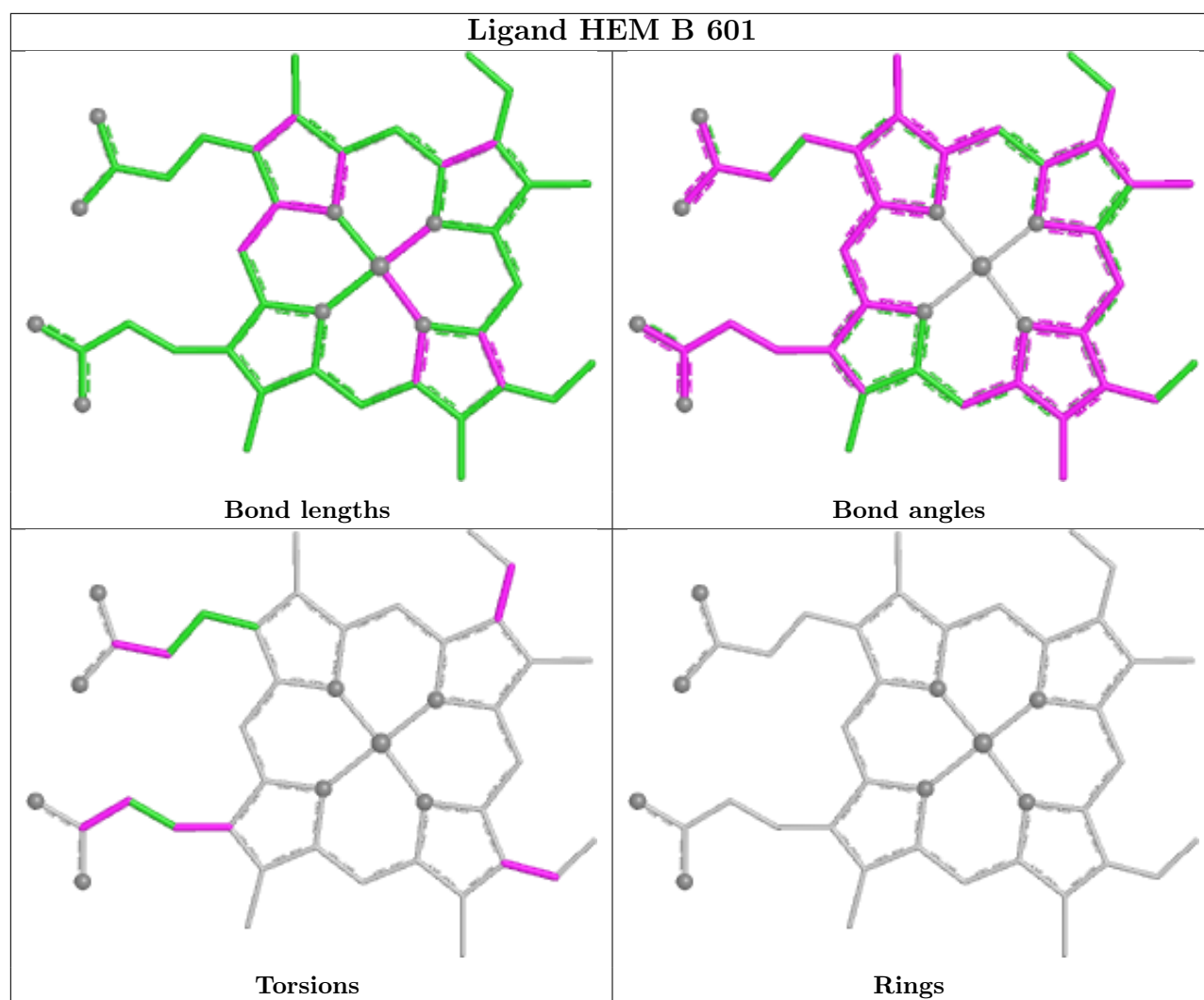
There are no ring outliers.

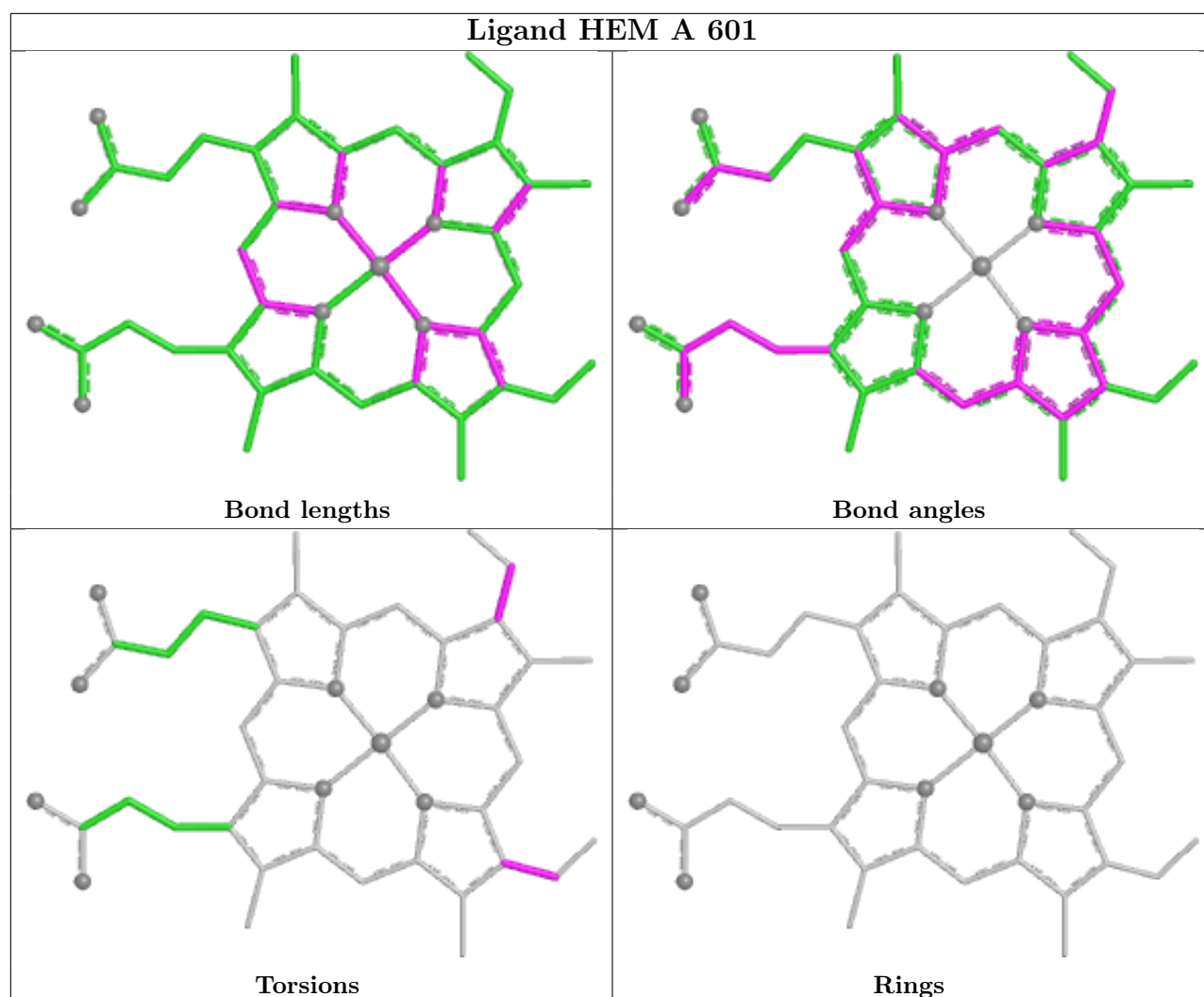
4 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	AER	4	0
3	B	602	AER	4	0
2	B	601	HEM	7	0
2	A	601	HEM	6	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	406/486 (83%)	0.03	10 (2%) 58 49	38, 90, 133, 160	4 (0%)
1	B	414/486 (85%)	-0.10	3 (0%) 84 80	48, 82, 125, 161	1 (0%)
All	All	820/972 (84%)	-0.04	13 (1%) 70 64	38, 86, 132, 161	5 (0%)

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	139	PHE	3.5
1	A	202	LEU	3.2
1	B	465	CYS	3.2
1	A	187	CYS	2.8
1	A	146	SER	2.7
1	B	242	CYS	2.6
1	A	453	PHE	2.6
1	A	145	GLY	2.3
1	A	449	GLU	2.2
1	A	208	TYR	2.2
1	A	134	LEU	2.2
1	B	453	PHE	2.1
1	A	276	GLY	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](#)

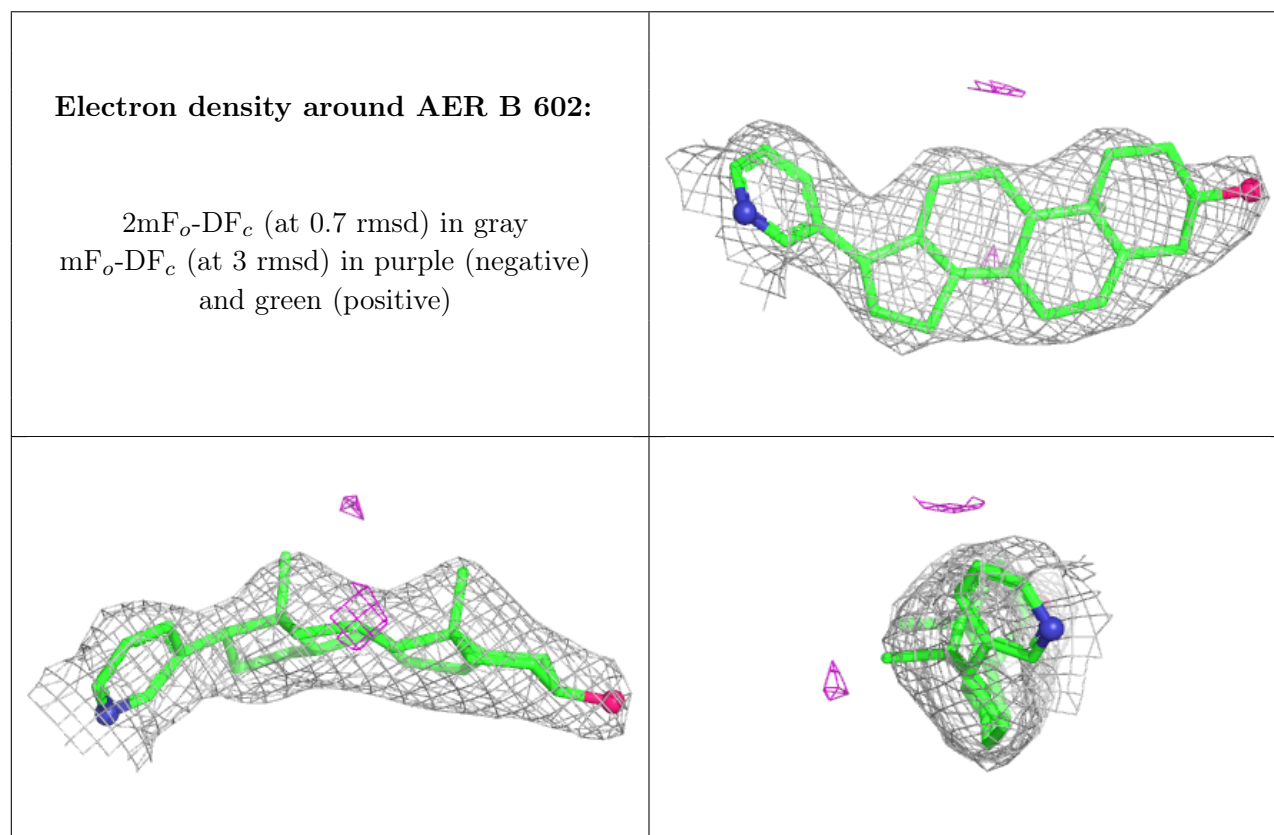
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

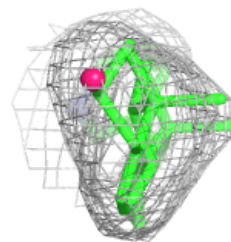
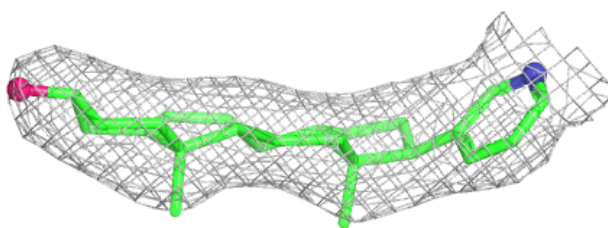
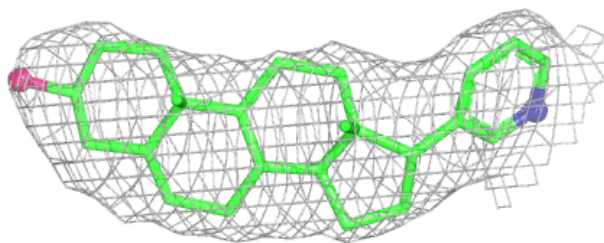
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	HG	B	603	1/1	0.81	0.07	152,152,152,152	1
3	AER	B	602	26/26	0.95	0.09	54,60,69,74	0
4	HG	A	604	1/1	0.95	0.07	116,116,116,116	1
4	HG	A	605	1/1	0.95	0.13	88,88,88,88	1
3	AER	A	602	26/26	0.95	0.09	55,66,72,75	0
4	HG	B	604	1/1	0.95	0.06	114,114,114,114	1
2	HEM	A	601	43/43	0.97	0.07	44,50,71,79	0
4	HG	A	603	1/1	0.97	0.04	126,126,126,126	1
2	HEM	B	601	43/43	0.98	0.06	38,44,53,58	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.



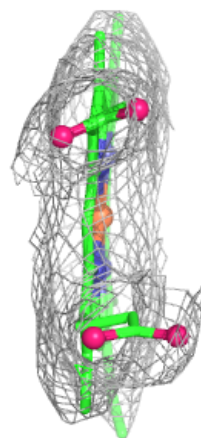
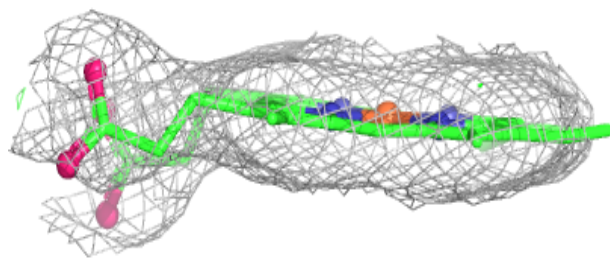
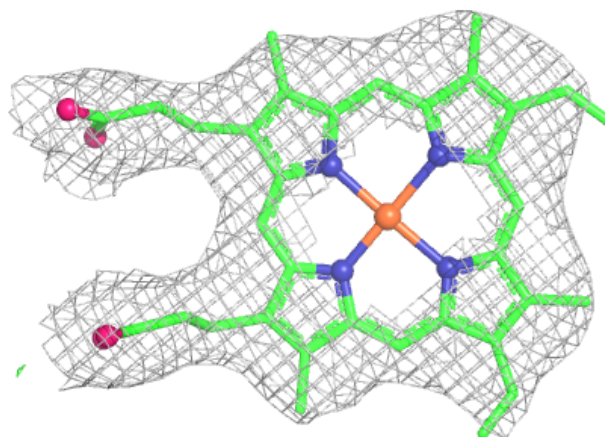
Electron density around AER A 602:

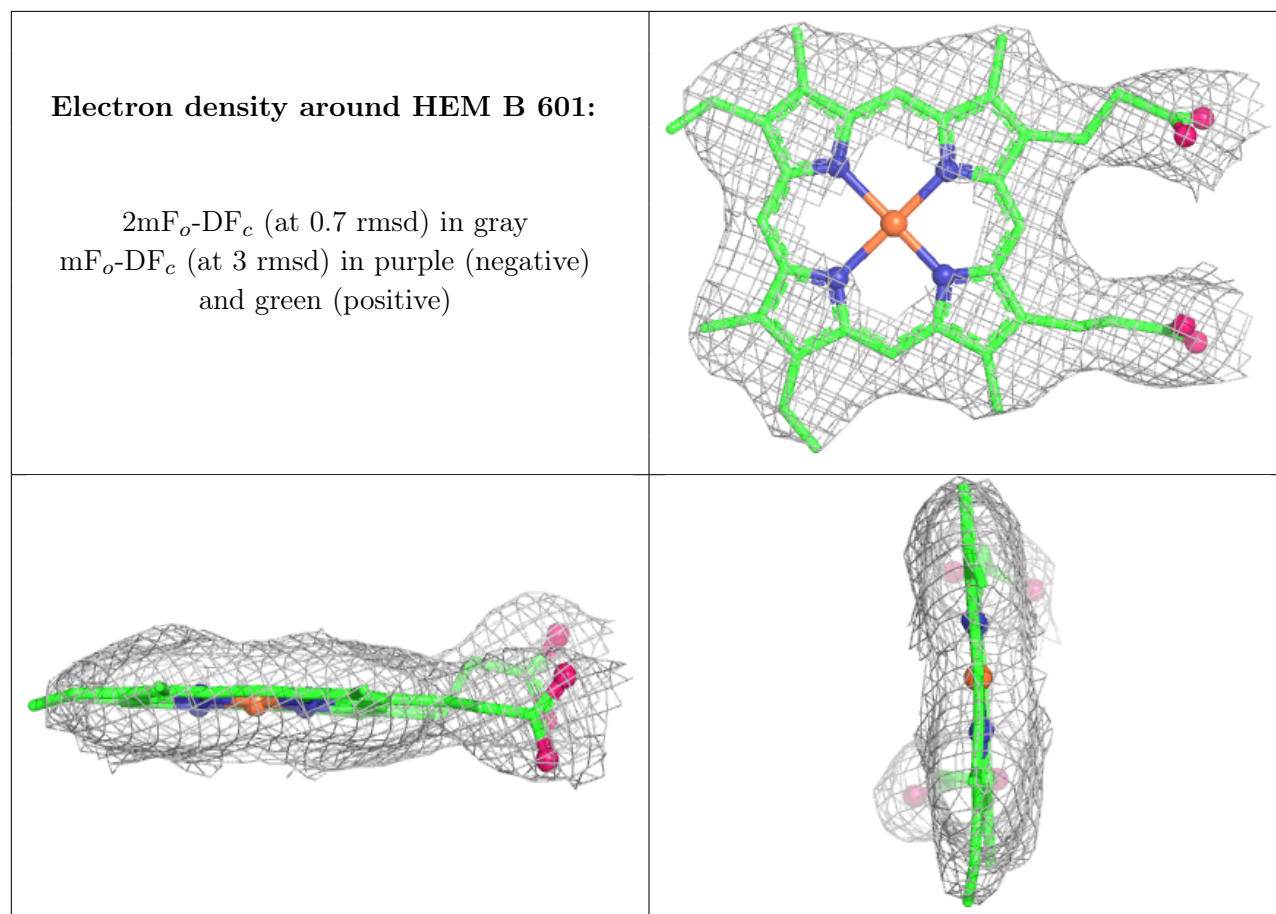
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 601:

$2mF_o-DF_c$ (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.