



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:10 PM UTC

PDB ID : 5TZ1 / pdb_00005tz1
Title : Crystal structure of sterol 14-alpha demethylase (CYP51) from *Candida albicans* in complex with the tetrazole-based antifungal drug candidate VT1161 (VT1)
Authors : Hargrove, T.; Wawrzak, Z.; Lepesheva, G.
Deposited on : 2016-11-21
Resolution : 2.00 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

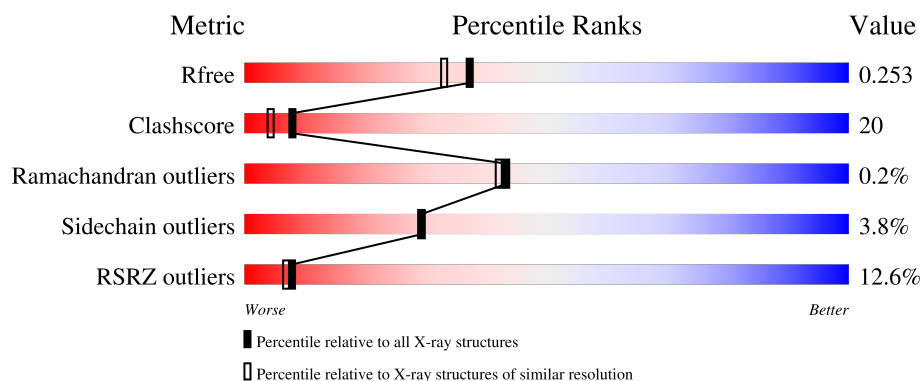
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	490	11% 77% 20% ..
1	B	490	14% 69% 25% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8333 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

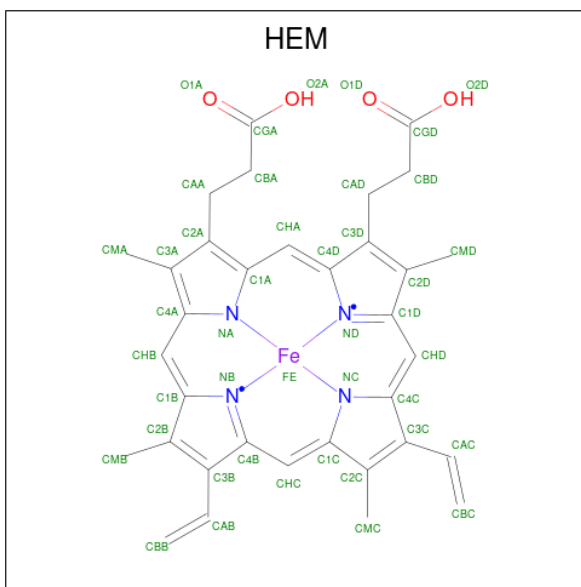
- Molecule 1 is a protein called Sterol 14-alpha demethylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	484	Total	C	N	O	S	0	0	0
			3928	2539	652	721	16			
1	B	484	Total	C	N	O	S	0	0	0
			3929	2539	652	722	16			

There are 22 discrepancies between the modelled and reference sequences:

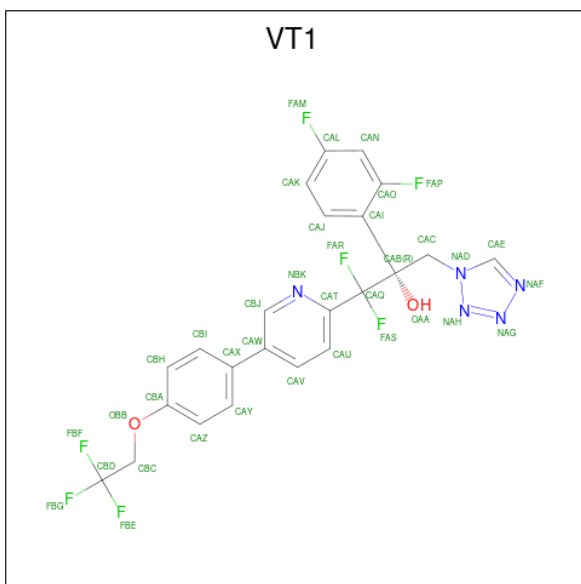
Chain	Residue	Modelled	Actual	Comment	Reference
A	43	MET	-	expression tag	UNP Q9P4W0
A	44	ALA	-	expression tag	UNP Q9P4W0
A	45	LYS	-	expression tag	UNP Q9P4W0
A	46	LYS	-	expression tag	UNP Q9P4W0
A	47	THR	-	expression tag	UNP Q9P4W0
A	48	PRO	ALA	engineered mutation	UNP Q9P4W0
A	263	LEU	SER	engineered mutation	UNP Q9P4W0
A	529	HIS	-	expression tag	UNP Q9P4W0
A	530	HIS	-	expression tag	UNP Q9P4W0
A	531	HIS	-	expression tag	UNP Q9P4W0
A	532	HIS	-	expression tag	UNP Q9P4W0
B	43	MET	-	expression tag	UNP Q9P4W0
B	44	ALA	-	expression tag	UNP Q9P4W0
B	45	LYS	-	expression tag	UNP Q9P4W0
B	46	LYS	-	expression tag	UNP Q9P4W0
B	47	THR	-	expression tag	UNP Q9P4W0
B	48	PRO	ALA	engineered mutation	UNP Q9P4W0
B	263	LEU	SER	engineered mutation	UNP Q9P4W0
B	529	HIS	-	expression tag	UNP Q9P4W0
B	530	HIS	-	expression tag	UNP Q9P4W0
B	531	HIS	-	expression tag	UNP Q9P4W0
B	532	HIS	-	expression tag	UNP Q9P4W0

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



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

- Molecule 3 is (R)-2-(2,4-Difluorophenyl)-1,1-difluoro-3-(1H-tetrazol-1-yl)-1-(5-(4-(2,2,2-trifluoroethoxy)phenyl)pyridin-2-yl)propan-2-ol (CCD ID: VT1) (formula: C₂₃H₁₆F₇N₅O₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			37	23	7	5	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	F	N	O	0	0
			37	23	7	5	2		

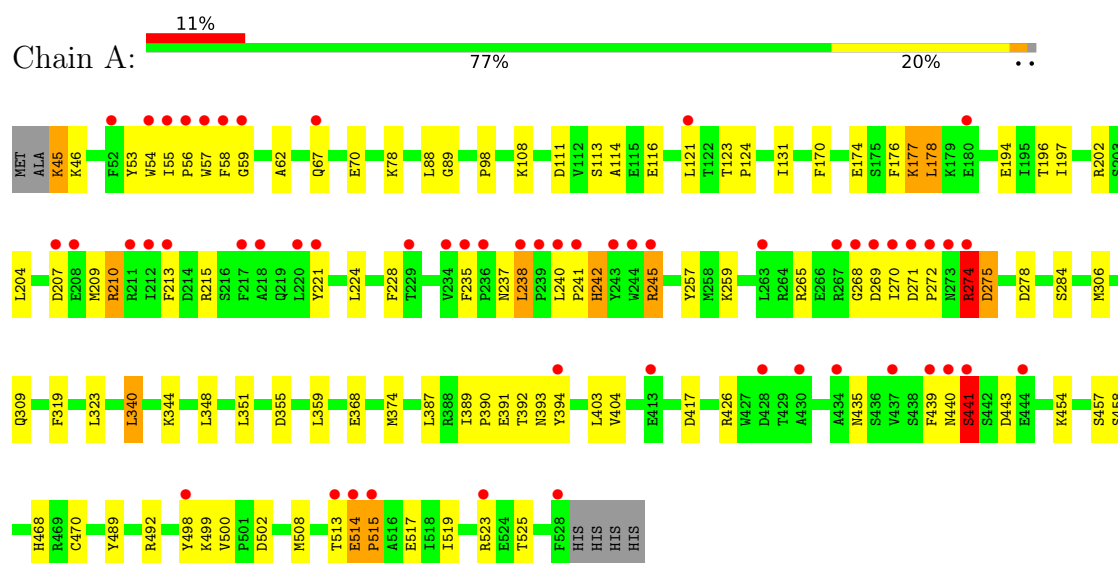
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	143	Total	O	0	0
			143	143		
4	B	173	Total	O	0	0
			173	173		

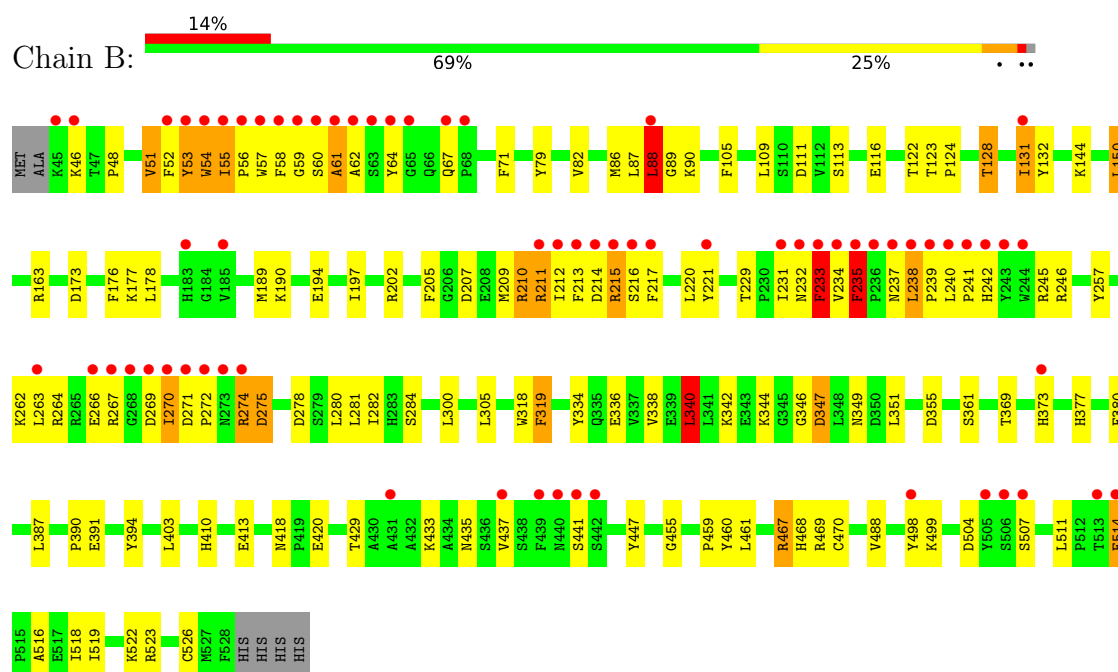
3 Residue-property plots

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: Sterol 14-alpha demethylase



• Molecule 1: Sterol 14-alpha demethylase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	177.64Å 71.44Å 79.19Å 90.00° 96.63° 90.00°	Depositor
Resolution (Å)	49.61 – 2.00 49.61 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.6 (49.61-2.00) 98.6 (49.61-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.218 , 0.224 0.231 , 0.253	Depositor DCC
R_{free} test set	3269 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	29.4	Xtriage
Anisotropy	0.259	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8333	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.79% 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, VT1

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	0.54	2/4039 (0.0%)	0.85	7/5474 (0.1%)
1	B	0.60	4/4040 (0.1%)	0.99	20/5475 (0.4%)
All	All	0.57	6/8079 (0.1%)	0.92	27/10949 (0.2%)

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	B	0	4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	518	ILE	C-O	-6.30	1.17	1.24
1	A	340	LEU	C-O	-5.60	1.17	1.24
1	B	235	PHE	C-O	-5.44	1.21	1.23
1	A	284	SER	C-O	-5.39	1.16	1.24
1	B	220	LEU	C-O	-5.39	1.18	1.24
1	B	340	LEU	C-O	-5.17	1.18	1.24

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	54	TRP	N-CA-C	16.55	129.02	111.14
1	B	51	VAL	N-CA-C	10.44	122.64	111.58
1	B	217	PHE	N-CA-C	-10.12	99.48	112.23
1	B	233	PHE	CA-C-N	9.46	132.71	122.11
1	B	233	PHE	C-N-CA	9.46	132.71	122.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	268	GLY	N-CA-C	8.88	121.74	111.36
1	B	61	ALA	N-CA-C	-8.36	103.10	113.38
1	B	55	ILE	CA-C-N	8.04	129.90	119.84
1	B	55	ILE	C-N-CA	8.04	129.90	119.84
1	B	514	GLU	N-CA-C	8.01	124.59	113.45
1	A	374	MET	N-CA-C	-7.74	99.81	109.64
1	B	275	ASP	N-CA-C	7.35	117.94	108.24
1	A	441	SER	N-CA-C	7.34	120.87	108.90
1	B	270	ILE	N-CA-C	-6.79	95.21	109.34
1	A	275	ASP	N-CA-C	6.31	116.40	108.45
1	B	88	LEU	CA-CB-CG	6.16	137.88	116.30
1	B	233	PHE	N-CA-C	-6.03	106.01	113.19
1	A	514	GLU	N-CA-C	5.90	122.85	109.81
1	B	380	PHE	N-CA-C	5.76	117.93	109.24
1	A	274	ARG	N-CA-C	5.74	117.34	111.14
1	B	238	LEU	CA-C-N	5.64	126.90	119.84
1	B	238	LEU	C-N-CA	5.64	126.90	119.84
1	B	150	LEU	O-C-N	-5.60	115.37	122.26
1	A	515	PRO	N-CA-C	-5.29	101.58	112.47
1	B	346	GLY	N-CA-C	5.23	118.74	111.10
1	B	216	SER	N-CA-C	-5.15	107.35	113.38
1	B	150	LEU	N-CA-C	5.13	118.85	112.59

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	150	LEU	Mainchain
1	B	233	PHE	Sidechain
1	B	319	PHE	Sidechain
1	B	347	ASP	Mainchain

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	3928	0	3861	128	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3929	0	3865	181	0
2	A	43	0	30	3	0
2	B	43	0	30	8	0
3	A	37	0	16	1	0
3	B	37	0	16	7	0
4	A	143	0	0	3	1
4	B	173	0	0	7	2
All	All	8333	0	7818	312	2

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 20.

All (312) 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:235:PHE:CD1	1:B:238:LEU:HD11	1.68	1.27
1:B:270:ILE:HD11	1:B:281:LEU:HD13	1.20	1.15
1:B:235:PHE:CG	1:B:238:LEU:HD11	1.81	1.14
1:B:55:ILE:O	1:B:58:PHE:HB3	1.54	1.07
1:B:52:PHE:CD1	1:B:54:TRP:HZ2	1.72	1.06
1:B:60:SER:O	1:B:64:TYR:HB3	1.59	1.02
1:A:270:ILE:N	1:A:274:ARG:HH12	1.60	1.00
1:A:274:ARG:HD3	1:A:278:ASP:OD2	1.61	1.00
1:B:87:LEU:HD23	1:B:88:LEU:HD12	1.42	0.99
1:B:211:ARG:HG3	1:B:211:ARG:HH11	1.25	0.97
1:B:52:PHE:CD1	1:B:54:TRP:CZ2	2.54	0.95
1:B:235:PHE:CE1	1:B:238:LEU:HD11	2.04	0.93
1:B:271:ASP:HB3	1:B:272:PRO:HD3	1.48	0.93
1:B:235:PHE:CG	1:B:238:LEU:CD1	2.53	0.92
1:B:235:PHE:CD2	1:B:238:LEU:CD1	2.54	0.91
1:B:270:ILE:HD11	1:B:281:LEU:CD1	2.00	0.90
1:B:207:ASP:HB3	1:B:211:ARG:HH12	1.36	0.90
1:A:271:ASP:CG	1:A:272:PRO:HD2	1.98	0.88
1:A:271:ASP:OD1	1:A:272:PRO:CD	2.22	0.87
1:A:269:ASP:C	1:A:274:ARG:HH12	1.84	0.84
1:B:269:ASP:OD1	1:B:271:ASP:HB2	1.77	0.84
1:B:240:LEU:HD13	1:B:242:HIS:ND1	1.92	0.84
1:B:131:ILE:HG22	1:B:132:TYR:N	1.93	0.83
2:A:601:HEM:HMB2	2:A:601:HEM:HBB2	1.60	0.83
1:A:271:ASP:O	1:A:274:ARG:HG3	1.78	0.83
1:A:54:TRP:CD1	1:B:54:TRP:HE3	1.96	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:ILE:HG22	1:B:132:TYR:H	1.43	0.82
1:A:269:ASP:HA	1:A:274:ARG:NH1	1.95	0.81
1:A:121:LEU:HD11	1:A:508:MET:CE	2.11	0.81
1:A:270:ILE:HG22	1:A:274:ARG:CZ	2.11	0.81
1:B:235:PHE:CE2	1:B:238:LEU:HD13	2.16	0.81
1:A:121:LEU:HD11	1:A:508:MET:HE1	1.63	0.80
1:A:269:ASP:CA	1:A:274:ARG:HH12	1.94	0.80
1:A:270:ILE:N	1:A:274:ARG:NH1	2.28	0.80
1:B:52:PHE:HD2	1:B:71:PHE:HE1	1.30	0.79
1:B:87:LEU:O	1:B:88:LEU:HD13	1.82	0.79
1:B:334:TYR:CE1	1:B:523:ARG:HD2	2.16	0.79
1:B:498:TYR:C	1:B:499:LYS:HD2	2.09	0.78
1:B:52:PHE:HD2	1:B:71:PHE:CE1	2.02	0.78
1:A:113:SER:OG	1:A:116:GLU:HG2	1.84	0.77
1:A:269:ASP:HA	1:A:274:ARG:HH12	1.48	0.77
1:B:344:LYS:HE3	1:B:355:ASP:OD2	1.84	0.77
1:A:271:ASP:OD1	1:A:272:PRO:HD2	1.82	0.77
1:B:52:PHE:HD1	1:B:54:TRP:HZ2	1.29	0.77
1:B:271:ASP:HB3	1:B:272:PRO:CD	2.14	0.77
1:B:51:VAL:O	1:B:52:PHE:CD1	2.37	0.77
1:A:271:ASP:OD1	1:A:272:PRO:HD3	1.84	0.76
1:A:265:ARG:NH1	4:A:701:HOH:O	1.96	0.76
1:B:235:PHE:CD2	1:B:238:LEU:HD11	2.20	0.76
1:B:232:ASN:HA	1:B:234:VAL:O	1.85	0.75
1:B:51:VAL:O	1:B:52:PHE:HD1	1.70	0.75
1:B:514:GLU:HB2	4:B:726:HOH:O	1.87	0.75
1:B:232:ASN:C	1:B:234:VAL:H	1.93	0.74
1:A:54:TRP:HD1	1:B:54:TRP:HE3	1.35	0.74
1:B:52:PHE:HD1	1:B:54:TRP:CZ2	2.03	0.73
1:B:57:TRP:C	1:B:88:LEU:HA	2.16	0.71
1:B:235:PHE:CE2	1:B:238:LEU:CD1	2.72	0.71
1:A:202:ARG:NE	1:A:207:ASP:OD1	2.23	0.71
1:A:111:ASP:HB3	1:A:387:LEU:HD21	1.73	0.70
1:B:57:TRP:O	1:B:88:LEU:CB	2.38	0.70
1:B:58:PHE:CE1	1:B:61:ALA:HB3	2.27	0.70
1:A:274:ARG:CD	1:A:274:ARG:H	2.04	0.70
1:B:270:ILE:CD1	1:B:281:LEU:HD13	2.12	0.70
1:B:57:TRP:HA	1:B:88:LEU:HA	1.75	0.69
1:A:176:PHE:HB3	1:A:178:LEU:HD13	1.75	0.69
1:A:54:TRP:NE1	1:B:54:TRP:O	2.25	0.69
1:A:274:ARG:HD3	1:A:274:ARG:H	1.58	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:271:ASP:O	1:A:274:ARG:CG	2.41	0.68
1:B:340:LEU:HD23	1:B:340:LEU:C	2.18	0.68
1:A:274:ARG:CD	1:A:278:ASP:OD2	2.39	0.68
1:B:274:ARG:HD2	1:B:278:ASP:OD2	1.94	0.68
1:A:270:ILE:H	1:A:274:ARG:NH1	1.91	0.67
1:B:207:ASP:HB3	1:B:211:ARG:NH1	2.08	0.67
1:B:237:ASN:O	1:B:238:LEU:HG	1.94	0.66
1:B:211:ARG:HH11	1:B:211:ARG:CG	2.06	0.66
1:B:334:TYR:CD1	1:B:523:ARG:HD2	2.31	0.66
1:B:202:ARG:HD3	1:B:210:ARG:HD3	1.77	0.65
1:B:87:LEU:O	1:B:88:LEU:CD1	2.45	0.65
2:B:601:HEM:HMC2	2:B:601:HEM:HBC2	1.79	0.64
1:A:241:PRO:O	1:A:245:ARG:HD3	1.96	0.64
1:B:87:LEU:CD2	1:B:88:LEU:HD12	2.24	0.64
1:A:389:ILE:O	1:A:392:THR:OG1	2.12	0.64
1:A:207:ASP:OD2	1:A:270:ILE:HD11	1.99	0.63
1:A:242:HIS:HA	1:A:245:ARG:NH1	2.13	0.63
1:B:90:LYS:HE2	1:B:233:PHE:CE1	2.33	0.63
1:A:215:ARG:CD	1:A:215:ARG:H	2.12	0.62
1:A:54:TRP:CD1	1:B:54:TRP:CE3	2.85	0.62
1:A:271:ASP:O	1:A:274:ARG:HD2	1.99	0.62
1:B:57:TRP:O	1:B:88:LEU:HB2	1.99	0.62
1:A:340:LEU:C	1:A:340:LEU:HD23	2.25	0.62
1:A:58:PHE:HE2	1:A:62:ALA:HB2	1.64	0.61
1:B:57:TRP:CA	1:B:88:LEU:HA	2.30	0.61
1:B:57:TRP:O	1:B:88:LEU:HB3	2.00	0.61
1:B:55:ILE:O	1:B:58:PHE:CB	2.41	0.61
1:B:52:PHE:CD2	1:B:71:PHE:HE1	2.17	0.61
1:A:403:LEU:C	1:A:403:LEU:HD23	2.26	0.61
1:B:240:LEU:CD1	1:B:242:HIS:ND1	2.63	0.61
1:B:270:ILE:HG21	1:B:282:ILE:HD11	1.82	0.61
1:B:235:PHE:CE1	1:B:238:LEU:CD1	2.82	0.60
1:A:108:LYS:HE3	1:A:111:ASP:OD2	2.01	0.60
1:B:88:LEU:HD22	1:B:89:GLY:H	1.67	0.60
2:B:601:HEM:HMB2	2:B:601:HEM:HBB2	1.82	0.60
1:A:441:SER:HB3	1:A:443:ASP:OD1	2.01	0.60
1:B:215:ARG:O	1:B:215:ARG:HG3	2.00	0.60
1:B:235:PHE:CD2	1:B:238:LEU:HD13	2.34	0.60
1:A:67:GLN:CG	1:A:70:GLU:HB2	2.32	0.59
1:A:213:PHE:CE2	1:A:221:TYR:HE2	2.21	0.59
1:B:211:ARG:HG3	1:B:211:ARG:NH1	2.07	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:602:VT1:HAC2	3:B:602:VT1:FAP	1.91	0.59
1:B:235:PHE:CZ	1:B:238:LEU:CD1	2.86	0.59
1:A:237:ASN:O	1:A:238:LEU:HD22	2.02	0.58
2:B:601:HEM:HBC2	2:B:601:HEM:CMC	2.34	0.58
1:A:54:TRP:HD1	1:B:54:TRP:CE3	2.19	0.58
1:B:278:ASP:O	1:B:282:ILE:HD12	2.04	0.58
1:B:232:ASN:C	1:B:234:VAL:N	2.59	0.58
1:B:231:ILE:O	1:B:234:VAL:O	2.22	0.58
1:B:270:ILE:CD1	1:B:281:LEU:CD1	2.77	0.57
1:B:111:ASP:HB3	1:B:387:LEU:HD21	1.85	0.57
1:B:271:ASP:CB	1:B:272:PRO:HD3	2.27	0.57
1:B:113:SER:OG	1:B:116:GLU:HG2	2.04	0.57
1:B:514:GLU:CB	4:B:726:HOH:O	2.50	0.57
1:B:272:PRO:HB2	1:B:274:ARG:HG2	1.86	0.57
1:A:270:ILE:H	1:A:274:ARG:HH12	1.42	0.57
1:B:526:CYS:N	4:B:703:HOH:O	2.38	0.57
1:B:235:PHE:CE1	1:B:238:LEU:HD21	2.40	0.56
1:B:51:VAL:HG11	1:B:79:TYR:CD2	2.40	0.56
1:A:67:GLN:HG2	1:A:70:GLU:CD	2.30	0.56
1:B:213:PHE:HD1	1:B:213:PHE:O	1.89	0.56
1:B:338:VAL:O	1:B:342:LYS:HG2	2.06	0.56
1:B:56:PRO:C	1:B:58:PHE:H	2.14	0.56
1:B:347:ASP:OD1	1:B:349:ASN:HB2	2.06	0.56
1:B:514:GLU:O	1:B:516:ALA:N	2.40	0.56
1:A:514:GLU:HB3	1:A:515:PRO:HD3	1.88	0.55
1:A:57:TRP:CH2	1:A:89:GLY:HA3	2.42	0.55
1:A:492:ARG:HG3	1:A:523:ARG:HH11	1.71	0.55
1:B:523:ARG:HB2	4:B:703:HOH:O	2.05	0.55
1:A:235:PHE:HB3	1:A:238:LEU:HD23	1.89	0.55
1:B:58:PHE:HE2	1:B:62:ALA:HB2	1.70	0.55
1:A:270:ILE:H	1:A:274:ARG:NH2	2.05	0.55
1:A:224:LEU:HD23	1:A:306:MET:HE1	1.89	0.55
1:B:429:THR:O	1:B:433:LYS:HG3	2.07	0.55
1:A:270:ILE:H	1:A:274:ARG:CZ	2.20	0.54
1:B:235:PHE:CZ	1:B:238:LEU:HD11	2.42	0.54
1:A:53:TYR:CD2	1:A:59:GLY:HA2	2.42	0.54
1:A:498:TYR:HH	1:B:413:GLU:CD	2.16	0.54
1:A:54:TRP:HE1	1:B:54:TRP:HA	1.73	0.54
1:B:122:THR:HB	4:B:842:HOH:O	2.08	0.54
1:B:53:TYR:HB3	1:B:86:MET:HB2	1.89	0.54
1:A:213:PHE:CE2	1:A:221:TYR:CE2	2.95	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:ILE:CG2	1:B:132:TYR:N	2.62	0.53
1:B:53:TYR:N	1:B:53:TYR:CD1	2.76	0.53
1:A:498:TYR:CD1	1:A:499:LYS:HD2	2.43	0.53
2:A:601:HEM:HBB2	2:A:601:HEM:CMB	2.33	0.53
1:A:271:ASP:O	1:A:274:ARG:CD	2.56	0.53
1:A:228:PHE:CE1	1:A:508:MET:HE2	2.44	0.53
1:B:280:LEU:O	1:B:284:SER:HB3	2.08	0.52
1:A:204:LEU:O	1:A:275:ASP:HB2	2.09	0.52
1:A:241:PRO:O	1:A:245:ARG:CD	2.57	0.52
1:B:205:PHE:O	4:B:701:HOH:O	2.19	0.52
1:A:228:PHE:CZ	1:A:508:MET:HE2	2.45	0.52
1:A:67:GLN:HG3	1:A:70:GLU:HB2	1.90	0.52
1:B:237:ASN:HD22	1:B:246:ARG:HH22	1.55	0.52
1:B:52:PHE:CD2	1:B:71:PHE:CE1	2.92	0.52
1:B:55:ILE:CG2	1:B:56:PRO:HD2	2.40	0.52
1:B:131:ILE:CG2	1:B:132:TYR:H	2.09	0.52
1:A:498:TYR:OH	1:B:413:GLU:OE1	2.28	0.52
1:A:242:HIS:HA	1:A:245:ARG:CZ	2.40	0.51
1:A:45:LYS:HG3	1:A:393:ASN:O	2.10	0.51
1:A:170:PHE:HB3	1:A:178:LEU:HD22	1.92	0.51
1:A:271:ASP:CG	1:A:272:PRO:CD	2.73	0.51
1:A:121:LEU:HD13	1:A:121:LEU:O	2.11	0.51
1:A:390:PRO:O	1:A:391:GLU:HB2	2.11	0.51
1:B:194:GLU:O	1:B:197:ILE:HG22	2.12	0.50
1:A:202:ARG:CD	1:A:210:ARG:HD2	2.41	0.50
1:B:488:VAL:O	1:B:523:ARG:NH1	2.44	0.50
1:A:499:LYS:HE3	1:B:413:GLU:OE2	2.12	0.50
1:A:523:ARG:NH1	4:A:710:HOH:O	2.45	0.50
1:A:209:MET:HE2	1:A:257:TYR:CD1	2.46	0.50
1:B:270:ILE:HG21	1:B:282:ILE:CD1	2.42	0.50
1:A:215:ARG:H	1:A:215:ARG:HD3	1.77	0.49
1:A:499:LYS:N	1:A:499:LYS:CD	2.74	0.49
1:B:123:THR:N	1:B:124:PRO:CD	2.74	0.49
1:A:470:CYS:HA	2:A:601:HEM:C4D	2.47	0.49
1:A:502:ASP:O	1:A:513:THR:HG23	2.12	0.49
1:A:121:LEU:HD13	1:A:121:LEU:C	2.37	0.49
1:B:262:LYS:O	1:B:266:GLU:HB2	2.13	0.49
1:B:468:HIS:HD2	2:B:601:HEM:O2D	1.95	0.49
1:A:194:GLU:O	1:A:197:ILE:HG22	2.12	0.49
1:A:270:ILE:H	1:A:274:ARG:HH22	1.60	0.49
1:B:338:VAL:HG13	1:B:342:LYS:CE	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:267:ARG:NH2	1:B:274:ARG:NH2	2.60	0.49
1:B:338:VAL:HG13	1:B:342:LYS:HE2	1.94	0.49
3:A:602:VT1:FAP	3:A:602:VT1:HAC2	2.02	0.48
1:B:232:ASN:CA	1:B:234:VAL:O	2.57	0.48
1:A:54:TRP:O	1:A:56:PRO:HD3	2.13	0.48
1:B:264:ARG:HH12	1:B:275:ASP:CG	2.21	0.48
1:A:245:ARG:HH11	1:A:245:ARG:HB2	1.77	0.48
1:B:435:ASN:O	1:B:459:PRO:HD2	2.14	0.48
1:A:270:ILE:CG2	1:A:274:ARG:HG2	2.44	0.48
1:B:235:PHE:CZ	1:B:238:LEU:HD13	2.49	0.48
1:A:319:PHE:CD1	1:A:319:PHE:C	2.92	0.47
1:A:351:LEU:C	1:A:351:LEU:HD23	2.39	0.47
1:B:336:GLU:OE2	1:B:361:SER:OG	2.30	0.47
1:A:174:GLU:HA	1:A:177:LYS:HD3	1.96	0.47
1:B:59:GLY:HA3	1:B:87:LEU:HA	1.95	0.47
1:A:196:THR:HG22	1:A:309:GLN:HG2	1.96	0.47
1:B:318:TRP:CE3	1:B:373:HIS:CE1	3.02	0.47
1:A:417:ASP:O	1:A:426:ARG:NH2	2.38	0.47
1:A:224:LEU:CD2	1:A:306:MET:HE1	2.45	0.47
1:B:58:PHE:CZ	1:B:61:ALA:HB3	2.50	0.47
1:A:57:TRP:CZ3	1:A:89:GLY:HA3	2.49	0.47
1:B:51:VAL:HG11	1:B:79:TYR:CE2	2.50	0.47
1:B:202:ARG:CD	1:B:210:ARG:HD3	2.42	0.47
1:A:123:THR:N	1:A:124:PRO:CD	2.78	0.46
1:A:492:ARG:HG3	1:A:523:ARG:NH1	2.29	0.46
1:B:144:LYS:HB3	1:B:144:LYS:HE3	1.84	0.46
1:B:498:TYR:O	1:B:499:LYS:HD2	2.14	0.46
1:B:213:PHE:O	1:B:213:PHE:CD1	2.67	0.46
1:A:319:PHE:O	1:A:323:LEU:HG	2.16	0.46
1:B:173:ASP:OD2	1:B:190:LYS:NZ	2.47	0.46
1:B:390:PRO:O	1:B:391:GLU:HB2	2.16	0.46
1:A:98:PRO:HB3	1:A:457:SER:O	2.15	0.46
1:B:238:LEU:HA	1:B:239:PRO:HD3	1.77	0.45
1:B:56:PRO:C	1:B:58:PHE:N	2.74	0.45
1:A:45:LYS:CD	1:A:393:ASN:HB3	2.46	0.45
1:B:48:PRO:HB3	1:B:82:VAL:HG12	1.98	0.45
1:A:209:MET:HE2	1:A:257:TYR:HD1	1.82	0.45
1:A:271:ASP:CB	1:A:272:PRO:HD2	2.46	0.45
1:A:340:LEU:HD23	1:A:340:LEU:O	2.17	0.45
1:B:274:ARG:H	1:B:274:ARG:HG3	1.49	0.45
1:B:410:HIS:CD2	1:B:460:TYR:HA	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:LEU:HD12	1:A:242:HIS:H	1.82	0.45
1:A:45:LYS:CG	1:A:393:ASN:HB3	2.46	0.45
1:A:368:GLU:OE2	1:A:426:ARG:HD3	2.16	0.45
1:A:240:LEU:CD1	1:A:242:HIS:HB3	2.48	0.45
1:B:189:MET:HE1	1:B:511:LEU:HD13	1.99	0.45
1:B:221:TYR:OH	1:B:305:LEU:HD23	2.17	0.45
1:B:90:LYS:CD	1:B:233:PHE:CZ	3.01	0.44
1:B:264:ARG:HD3	1:B:278:ASP:OD1	2.17	0.44
1:B:274:ARG:CD	1:B:278:ASP:OD2	2.63	0.44
1:A:45:LYS:HD3	1:A:393:ASN:HB3	1.99	0.44
1:B:469:ARG:HG2	1:B:470:CYS:N	2.31	0.44
1:A:435:ASN:O	1:A:458:SER:OG	2.35	0.44
1:B:55:ILE:HG23	1:B:56:PRO:HD2	2.00	0.44
1:B:235:PHE:CE2	1:B:238:LEU:HD11	2.48	0.44
1:B:319:PHE:HB3	1:B:369:THR:OG1	2.17	0.44
1:B:499:LYS:HD2	1:B:499:LYS:N	2.33	0.44
1:A:67:GLN:HG2	1:A:70:GLU:CG	2.47	0.44
1:B:377:HIS:NE2	3:B:602:VT1:OBB	2.51	0.44
3:B:602:VT1:FAP	3:B:602:VT1:CAC	2.55	0.44
1:A:202:ARG:HD3	1:A:210:ARG:HD2	1.98	0.44
2:B:601:HEM:HBB2	2:B:601:HEM:CMB	2.48	0.44
1:B:109:LEU:HD23	1:B:109:LEU:HA	1.87	0.44
1:A:78:LYS:NZ	4:A:715:HOH:O	2.50	0.44
1:B:344:LYS:HE3	1:B:355:ASP:CG	2.42	0.44
1:A:113:SER:CB	1:A:116:GLU:HG2	2.47	0.44
1:B:111:ASP:CB	1:B:387:LEU:HD21	2.48	0.43
1:A:340:LEU:HD13	1:A:359:LEU:HD21	1.99	0.43
1:B:128:THR:HG22	4:B:833:HOH:O	2.18	0.43
1:B:377:HIS:CE1	3:B:602:VT1:OBB	2.71	0.43
1:B:470:CYS:HA	2:B:601:HEM:C4D	2.53	0.43
1:B:64:TYR:CE2	1:B:87:LEU:HD12	2.53	0.43
1:A:237:ASN:C	1:A:238:LEU:HD22	2.43	0.43
1:B:55:ILE:CG2	1:B:56:PRO:CD	2.97	0.43
1:A:46:LYS:O	1:A:394:TYR:HA	2.18	0.43
1:A:498:TYR:CZ	1:B:418:ASN:N	2.86	0.43
1:B:403:LEU:HD23	1:B:403:LEU:C	2.44	0.43
1:B:519:ILE:N	1:B:519:ILE:HD12	2.34	0.43
2:B:601:HEM:HMC2	2:B:601:HEM:CBC	2.47	0.43
1:B:334:TYR:CE1	1:B:523:ARG:CD	2.95	0.43
1:B:270:ILE:CG2	1:B:282:ILE:HD11	2.49	0.42
1:A:499:LYS:H	1:A:499:LYS:HD3	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:105:PHE:CE2	1:B:461:LEU:HD23	2.54	0.42
1:A:241:PRO:C	1:A:245:ARG:CD	2.92	0.42
1:B:178:LEU:O	1:B:522:LYS:NZ	2.37	0.42
1:B:215:ARG:HH11	1:B:215:ARG:CG	2.32	0.42
3:B:602:VT1:HBC2	3:B:602:VT1:HAZ	1.64	0.42
1:B:340:LEU:C	1:B:340:LEU:CD2	2.88	0.42
1:B:209:MET:HE2	1:B:257:TYR:CD1	2.55	0.42
1:B:241:PRO:O	1:B:242:HIS:C	2.62	0.42
1:B:46:LYS:O	1:B:394:TYR:HA	2.20	0.42
1:B:441:SER:HB2	1:B:455:GLY:CA	2.50	0.42
1:B:504:ASP:OD1	1:B:504:ASP:C	2.62	0.42
1:B:64:TYR:HE2	1:B:87:LEU:CD1	2.33	0.42
1:B:163:ARG:NH1	1:B:351:LEU:HB3	2.35	0.42
1:A:57:TRP:CZ2	1:A:88:LEU:O	2.73	0.42
1:A:348:LEU:HB2	1:A:489:TYR:CE2	2.55	0.42
1:B:51:VAL:C	1:B:52:PHE:CD1	2.97	0.41
1:B:176:PHE:O	1:B:177:LYS:C	2.62	0.41
1:A:519:ILE:HD12	1:A:519:ILE:N	2.35	0.41
1:B:55:ILE:HG23	1:B:56:PRO:CD	2.50	0.41
1:B:342:LYS:HA	1:B:342:LYS:HD3	1.77	0.41
2:B:601:HEM:HBD1	3:B:602:VT1:HAK	2.01	0.41
1:A:523:ARG:C	1:A:525:THR:N	2.77	0.41
1:B:113:SER:CB	1:B:116:GLU:HG2	2.51	0.41
1:A:319:PHE:C	1:A:319:PHE:HD1	2.28	0.41
1:A:259:LYS:HE2	1:A:259:LYS:HB2	1.84	0.41
1:B:116:GLU:OE1	1:B:116:GLU:HA	2.21	0.41
1:B:231:ILE:O	1:B:234:VAL:CG2	2.68	0.41
1:A:344:LYS:HE3	1:A:355:ASP:OD2	2.20	0.41
1:B:241:PRO:O	1:B:245:ARG:HG2	2.21	0.41
1:B:267:ARG:HH22	1:B:274:ARG:NH2	2.19	0.41
1:B:369:THR:O	1:B:373:HIS:N	2.46	0.41
1:B:447:TYR:CD2	1:B:467:ARG:HA	2.56	0.41
1:B:318:TRP:CZ3	1:B:373:HIS:CE1	3.09	0.41
1:A:340:LEU:C	1:A:340:LEU:CD2	2.93	0.41
1:A:403:LEU:HD23	1:A:404:VAL:N	2.35	0.41
1:A:443:ASP:OD2	1:A:454:LYS:HD3	2.21	0.41
3:B:602:VT1:NBK	3:B:602:VT1:HAC1	2.35	0.41
1:A:114:ALA:HB3	1:A:468:HIS:CE1	2.56	0.40
1:A:403:LEU:C	1:A:403:LEU:CD2	2.93	0.40
1:B:52:PHE:HA	1:B:54:TRP:CZ2	2.57	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:859:HOH:O	4:B:859:HOH:O[2_656]	2.06	0.14
4:A:829:HOH:O	4:B:770:HOH:O[1_564]	2.11	0.09

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	482/490 (98%)	467 (97%)	14 (3%)	1 (0%)	43	42
1	B	482/490 (98%)	464 (96%)	17 (4%)	1 (0%)	43	42
All	All	964/980 (98%)	931 (97%)	31 (3%)	2 (0%)	43	42

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	131	ILE
1	B	131	ILE

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	430/436 (99%)	416 (97%)	14 (3%)	33	34
1	B	431/436 (99%)	412 (96%)	19 (4%)	25	24
All	All	861/872 (99%)	828 (96%)	33 (4%)	29	29

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	LYS
1	A	55	ILE
1	A	177	LYS
1	A	178	LEU
1	A	210	ARG
1	A	238	LEU
1	A	242	HIS
1	A	245	ARG
1	A	274	ARG
1	A	439	PHE
1	A	440	ASN
1	A	441	SER
1	A	500	VAL
1	A	517	GLU
1	B	53	TYR
1	B	67	GLN
1	B	88	LEU
1	B	128	THR
1	B	210	ARG
1	B	211	ARG
1	B	212	ILE
1	B	214	ASP
1	B	215	ARG
1	B	229	THR
1	B	235	PHE
1	B	263	LEU
1	B	274	ARG
1	B	300	LEU
1	B	340	LEU
1	B	420	GLU
1	B	437	VAL
1	B	467	ARG
1	B	507	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	237	ASN
1	A	283	HIS
1	B	67	GLN
1	B	237	ASN

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Mol	Chain	Res	Type
1	B	295	GLN
1	B	363	ASN
1	B	400	HIS
1	B	440	ASN
1	B	468	HIS
1	B	474	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 ⓘ

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	HEM	A	601	1,3	50,50,50	2.05	18 (36%)	67,82,82	1.45	11 (16%)
3	VT1	A	602	2	37,40,40	3.60	6 (16%)	53,60,60	2.09	14 (26%)
3	VT1	B	602	2	37,40,40	3.61	6 (16%)	53,60,60	2.06	14 (26%)
2	HEM	B	601	1,3	50,50,50	2.12	17 (34%)	67,82,82	1.55	15 (22%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.
'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	601	1,3	-	1/14/54/54	-
3	VT1	A	602	2	-	5/33/36/36	0/4/4/4
3	VT1	B	602	2	-	5/33/36/36	0/4/4/4
2	HEM	B	601	1,3	-	2/14/54/54	-

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	602	VT1	CAB-CAI	-13.88	1.39	1.52
3	A	602	VT1	CAB-CAI	-13.78	1.39	1.52
3	A	602	VT1	NAD-NAH	-11.85	1.15	1.34
3	B	602	VT1	NAD-NAH	-11.73	1.16	1.34
3	B	602	VT1	NAF-NAG	-9.31	1.15	1.35
3	A	602	VT1	NAF-NAG	-9.26	1.15	1.35
3	B	602	VT1	NAH-NAG	-5.09	1.19	1.30
3	A	602	VT1	NAH-NAG	-5.07	1.19	1.30
2	B	601	HEM	FE-NB	4.94	2.10	1.94
2	A	601	HEM	FE-NB	4.92	2.10	1.94
2	A	601	HEM	C4D-ND	-4.90	1.31	1.40
2	B	601	HEM	C4D-ND	-4.61	1.32	1.40
2	B	601	HEM	C1B-NB	-4.45	1.32	1.40
2	B	601	HEM	C4B-NB	-4.41	1.30	1.38
3	B	602	VT1	CAW-CAX	-3.96	1.39	1.49
3	A	602	VT1	CAW-CAX	-3.94	1.39	1.49
2	A	601	HEM	C4B-NB	-3.87	1.31	1.38
2	A	601	HEM	C1B-NB	-3.83	1.33	1.40
2	B	601	HEM	C1D-ND	-3.71	1.31	1.38
2	B	601	HEM	FE-NC	3.71	2.07	1.95
2	A	601	HEM	C1C-C2C	-3.66	1.38	1.45
2	A	601	HEM	C1C-NC	-3.54	1.32	1.39
2	B	601	HEM	C1C-NC	-3.51	1.33	1.39
2	B	601	HEM	C4C-NC	-3.31	1.33	1.39
2	B	601	HEM	C1C-C2C	-3.29	1.38	1.45
2	A	601	HEM	FE-NC	3.27	2.06	1.95
2	A	601	HEM	C4C-NC	-3.26	1.33	1.39
2	A	601	HEM	O2A-CGA	-3.24	1.20	1.30
2	B	601	HEM	O2D-CGD	-3.21	1.20	1.30
3	A	602	VT1	CBJ-NBK	3.05	1.40	1.34
3	B	602	VT1	CBJ-NBK	3.05	1.40	1.34
2	B	601	HEM	O2A-CGA	-3.02	1.20	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	HEM	FE-ND	-2.89	1.86	1.94
2	A	601	HEM	O2D-CGD	-2.89	1.21	1.30
2	A	601	HEM	C2A-C3A	-2.62	1.32	1.38
2	A	601	HEM	C1D-ND	-2.57	1.33	1.38
2	B	601	HEM	C4A-NA	-2.46	1.35	1.39
2	A	601	HEM	FE-ND	-2.44	1.87	1.94
2	B	601	HEM	C2A-C3A	-2.37	1.32	1.38
2	B	601	HEM	C3C-C4C	-2.36	1.41	1.46
2	B	601	HEM	C3D-C2D	-2.34	1.31	1.36
2	A	601	HEM	C1A-NA	-2.25	1.35	1.39
2	B	601	HEM	C1A-C2A	-2.18	1.40	1.44
2	A	601	HEM	C4A-NA	-2.04	1.35	1.39
2	A	601	HEM	C1A-C2A	-2.01	1.40	1.44
2	A	601	HEM	C3C-C4C	-2.00	1.42	1.46
2	A	601	HEM	C3D-C2D	-2.00	1.32	1.36

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	VT1	CAB-CAC-NAD	-6.73	103.25	112.33
3	B	602	VT1	CAB-CAC-NAD	-6.33	103.78	112.33
3	A	602	VT1	NAD-CAE-NAF	-5.68	103.35	109.46
3	B	602	VT1	NAD-CAE-NAF	-5.63	103.41	109.46
3	B	602	VT1	NAD-NAH-NAG	5.22	112.88	106.42
3	A	602	VT1	NAD-NAH-NAG	5.08	112.71	106.42
2	B	601	HEM	CHC-C4B-NB	4.48	129.25	124.42
2	B	601	HEM	C1B-NB-C4B	3.98	109.92	105.21
3	A	602	VT1	CAC-NAD-NAH	3.87	126.31	120.43
3	B	602	VT1	CAE-NAD-NAH	-3.72	105.24	107.97
2	A	601	HEM	CHC-C4B-NB	3.72	128.43	124.42
2	B	601	HEM	CHD-C1D-ND	3.42	128.10	124.42
2	B	601	HEM	CHA-C4D-ND	3.41	128.58	124.37
2	A	601	HEM	O2D-CGD-O1D	-3.39	114.62	123.33
3	A	602	VT1	CAE-NAD-NAH	-3.37	105.50	107.97
3	A	602	VT1	CAO-CAN-CAL	3.33	120.22	116.67
3	B	602	VT1	CAN-CAO-CAI	-3.30	120.30	123.88
3	A	602	VT1	CBJ-NBK-CAT	3.29	121.12	117.60
3	A	602	VT1	CAN-CAO-CAI	-3.27	120.32	123.88
3	B	602	VT1	CAO-CAN-CAL	3.25	120.14	116.67
3	B	602	VT1	CAC-NAD-NAH	3.24	125.34	120.43
2	A	601	HEM	C1B-NB-C4B	3.23	109.03	105.21
3	B	602	VT1	CBJ-NBK-CAT	3.23	121.05	117.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	HEM	CHA-C4D-ND	3.07	128.17	124.37
2	B	601	HEM	O2A-CGA-CBA	2.89	123.13	114.00
3	A	602	VT1	CAU-CAT-NBK	-2.88	118.85	122.33
2	A	601	HEM	O2A-CGA-CBA	2.84	122.96	114.00
2	A	601	HEM	O2D-CGD-CBD	2.80	122.84	114.00
2	A	601	HEM	CHA-C4D-C3D	-2.76	120.14	125.23
3	A	602	VT1	CAK-CAL-CAN	-2.71	119.65	123.23
3	B	602	VT1	OBB-CBC-CBD	-2.67	101.76	108.00
2	B	601	HEM	C4C-NC-C1C	2.66	110.16	105.82
3	B	602	VT1	CAK-CAL-CAN	-2.65	119.74	123.23
3	B	602	VT1	CAU-CAT-NBK	-2.62	119.17	122.33
2	B	601	HEM	CHD-C1D-C2D	-2.60	120.93	125.03
3	B	602	VT1	CAB-CAI-CAO	-2.58	120.54	122.83
3	A	602	VT1	OBB-CBC-CBD	-2.58	101.97	108.00
2	B	601	HEM	O2D-CGD-O1D	-2.57	116.71	123.33
3	B	602	VT1	CAW-CBJ-NBK	-2.57	120.20	124.28
2	B	601	HEM	CHA-C4D-C3D	-2.56	120.51	125.23
2	A	601	HEM	CHD-C1D-ND	2.48	127.09	124.42
2	B	601	HEM	O2D-CGD-CBD	2.45	121.74	114.00
2	A	601	HEM	CBD-CAD-C3D	2.40	119.17	112.53
3	A	602	VT1	CAW-CBJ-NBK	-2.38	120.50	124.28
2	A	601	HEM	O2A-CGA-O1A	-2.29	117.43	123.33
2	B	601	HEM	O2A-CGA-O1A	-2.29	117.45	123.33
2	A	601	HEM	C4C-NC-C1C	2.24	109.48	105.82
3	B	602	VT1	CAJ-CAI-CAO	2.20	119.38	116.28
2	B	601	HEM	CHB-C1B-NB	2.17	127.05	124.37
3	A	602	VT1	CAJ-CAI-CAO	2.10	119.24	116.28
2	B	601	HEM	C3B-C4B-NB	-2.06	107.99	109.47
3	A	602	VT1	CBC-OBB-CBA	-2.05	111.61	117.80
2	B	601	HEM	C1A-CHA-C4D	-2.02	121.50	126.25
2	B	601	HEM	CBD-CAD-C3D	2.01	118.09	112.53

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	602	VT1	CAB-CAC-NAD-NAH
3	B	602	VT1	FAR-CAQ-CAT-CAU
3	B	602	VT1	FAS-CAQ-CAT-CAU
3	A	602	VT1	CAB-CAC-NAD-NAH
3	A	602	VT1	CAZ-CBA-OBB-CBC
3	A	602	VT1	CBH-CBA-OBB-CBC

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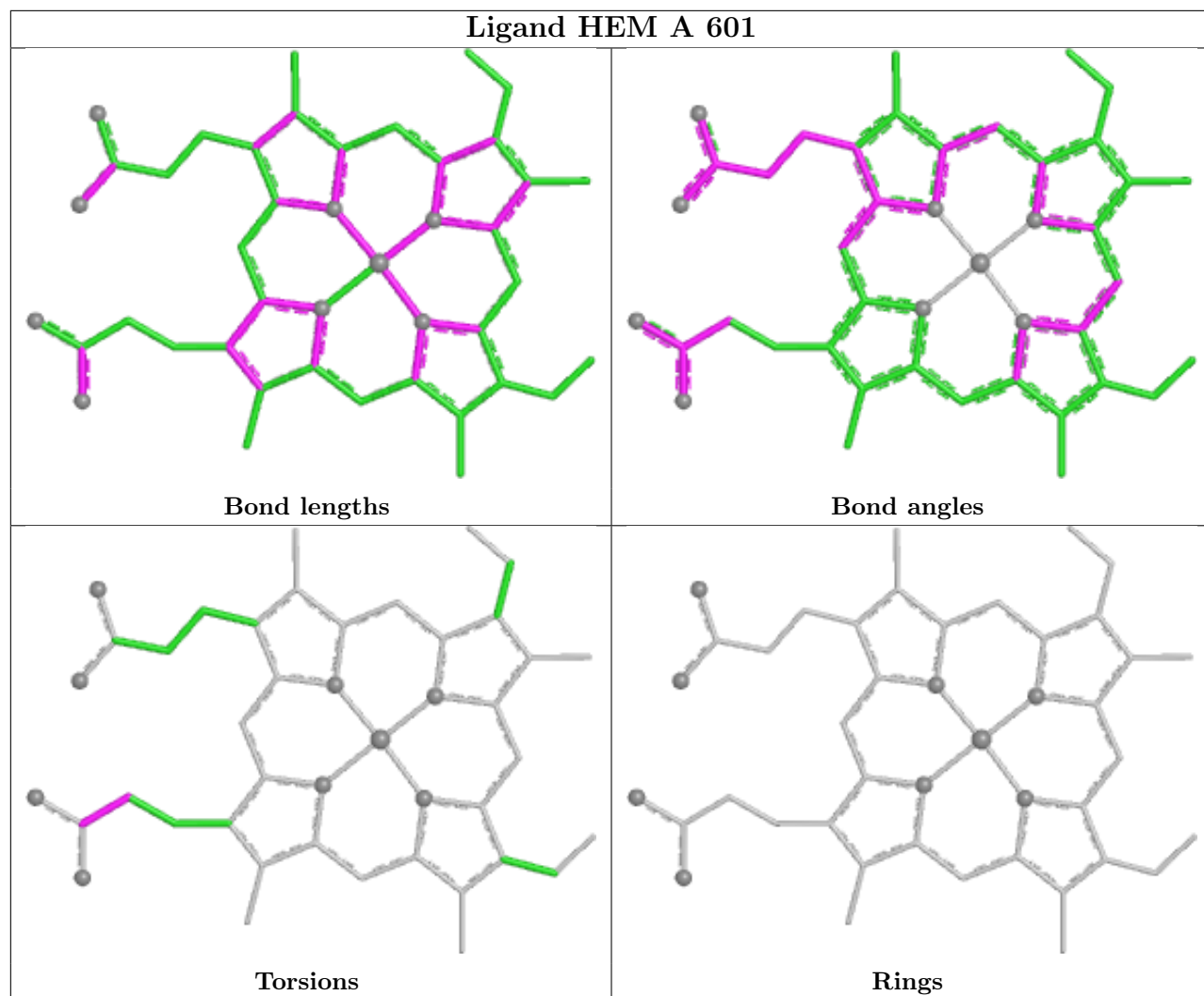
Mol	Chain	Res	Type	Atoms
3	B	602	VT1	CAB-CAQ-CAT-CAU
2	B	601	HEM	CAA-CBA-CGA-O2A
2	A	601	HEM	CAA-CBA-CGA-O2A
3	A	602	VT1	OBB-CBC-CBD-FBE
3	B	602	VT1	CAZ-CBA-OBB-CBC
3	A	602	VT1	OBB-CBC-CBD-FBF
2	B	601	HEM	CAA-CBA-CGA-O1A

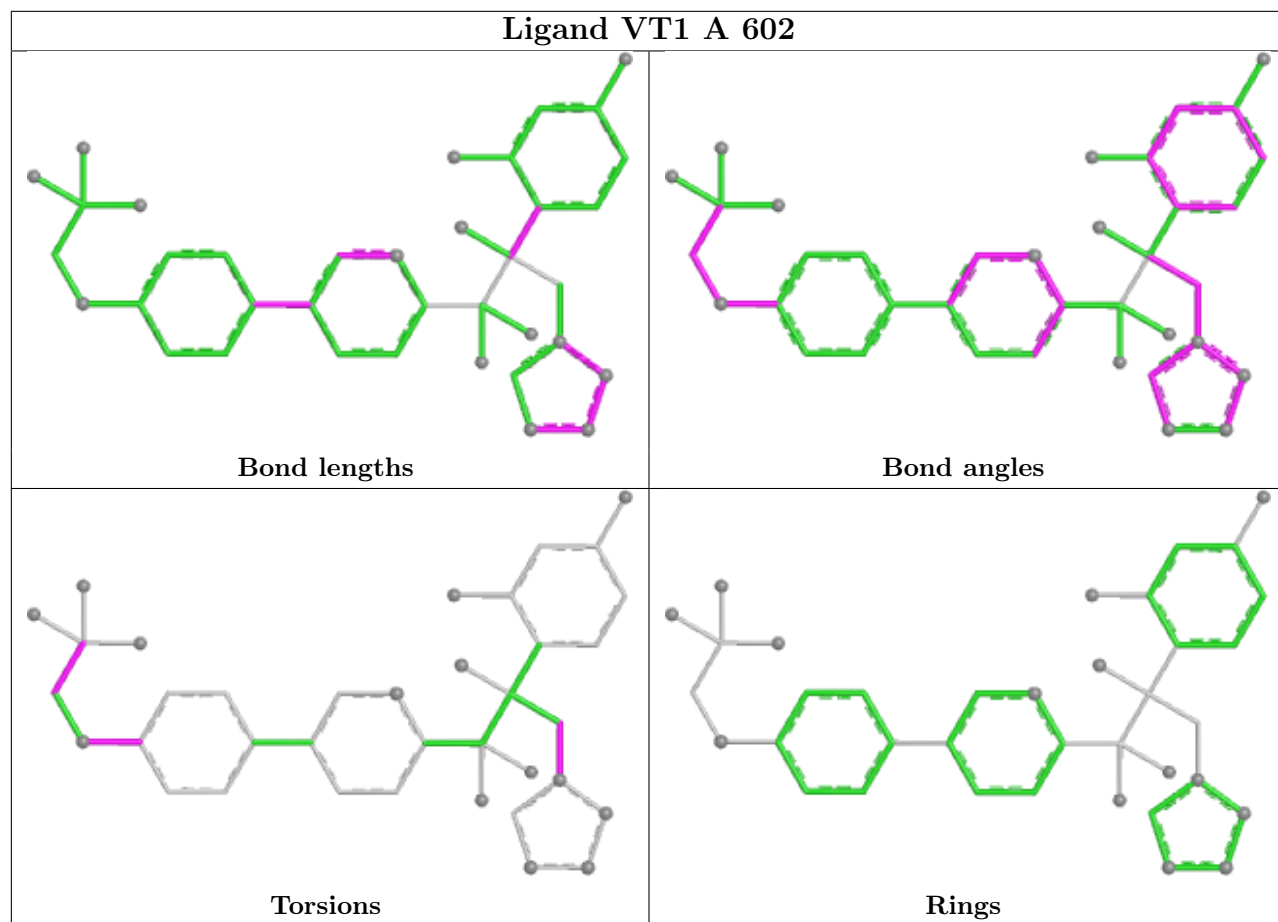
There are no ring outliers.

4 monomers are involved in 18 short contacts:

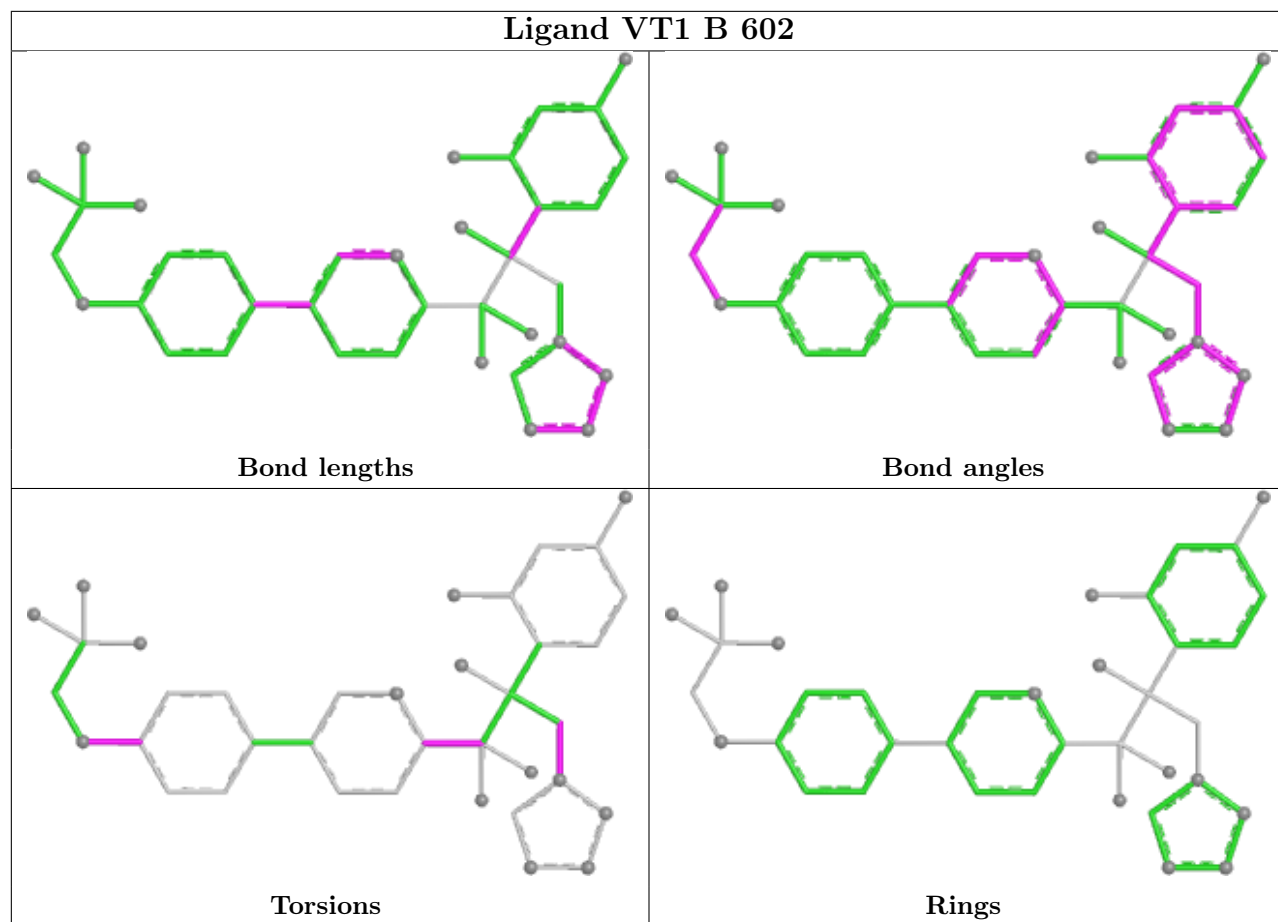
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	HEM	3	0
3	A	602	VT1	1	0
3	B	602	VT1	7	0
2	B	601	HEM	8	0

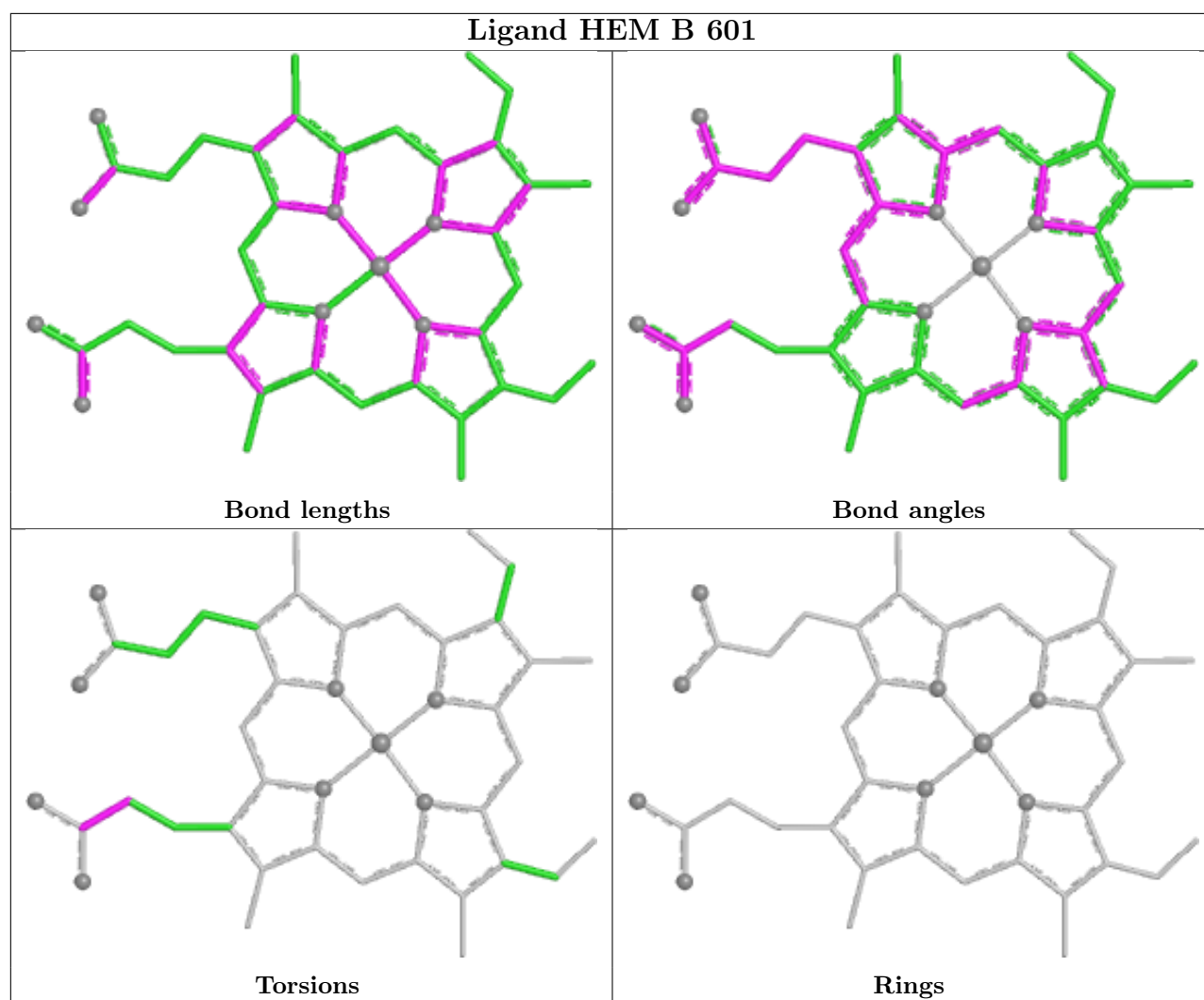
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 VT1 B 602





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	484/490 (98%)	0.78	55 (11%) 10 9	21, 35, 75, 123	0
1	B	484/490 (98%)	0.91	67 (13%) 6 6	21, 34, 94, 255	0
All	All	968/980 (98%)	0.84	122 (12%) 8 7	21, 34, 83, 255	0

All (122) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	54	TRP	8.4
1	B	55	ILE	7.4
1	B	57	TRP	7.2
1	B	272	PRO	7.1
1	A	514	GLU	6.7
1	B	60	SER	6.6
1	B	52	PHE	6.2
1	A	272	PRO	6.0
1	A	54	TRP	5.9
1	B	234	VAL	5.8
1	B	53	TYR	5.8
1	B	64	TYR	5.7
1	B	56	PRO	5.4
1	B	240	LEU	5.4
1	B	270	ILE	5.3
1	A	498	TYR	5.1
1	B	58	PHE	5.1
1	A	273	ASN	5.0
1	B	238	LEU	5.0
1	B	233	PHE	4.9
1	B	439	PHE	4.9
1	A	270	ILE	4.8
1	B	268	GLY	4.8
1	B	239	PRO	4.7

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Mol	Chain	Res	Type	RSRZ
1	B	441	SER	4.7
1	A	217	PHE	4.5
1	B	63	SER	4.5
1	A	440	ASN	4.5
1	B	244	TRP	4.3
1	B	241	PRO	4.3
1	B	267	ARG	4.2
1	B	65	GLY	4.2
1	A	58	PHE	4.2
1	B	213	PHE	4.1
1	B	216	SER	4.1
1	B	274	ARG	4.0
1	B	514	GLU	4.0
1	A	268	GLY	3.9
1	B	212	ILE	3.8
1	B	271	ASP	3.8
1	A	221	TYR	3.7
1	A	439	PHE	3.7
1	A	213	PHE	3.7
1	B	236	PRO	3.6
1	A	271	ASP	3.6
1	A	180	GLU	3.5
1	B	62	ALA	3.5
1	A	269	ASP	3.5
1	B	61	ALA	3.4
1	B	217	PHE	3.4
1	B	266	GLU	3.3
1	B	269	ASP	3.3
1	B	498	TYR	3.3
1	B	237	ASN	3.3
1	B	214	ASP	3.3
1	A	240	LEU	3.2
1	B	235	PHE	3.2
1	A	274	ARG	3.1
1	B	211	ARG	3.1
1	B	221	TYR	3.1
1	A	57	TRP	3.1
1	B	131	ILE	3.1
1	A	55	ILE	3.1
1	A	513	THR	3.0
1	B	273	ASN	3.0
1	B	243	TYR	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	45	LYS	3.0
1	B	59	GLY	2.9
1	B	513	THR	2.9
1	A	67	GLN	2.9
1	B	242	HIS	2.9
1	A	212	ILE	2.8
1	A	220	LEU	2.8
1	A	244	TRP	2.8
1	A	239	PRO	2.7
1	A	437	VAL	2.7
1	B	88	LEU	2.6
1	A	243	TYR	2.6
1	B	183	HIS	2.6
1	A	413	GLU	2.6
1	B	507	SER	2.6
1	A	235	PHE	2.5
1	A	59	GLY	2.5
1	B	215	ARG	2.5
1	A	238	LEU	2.5
1	B	506	SER	2.5
1	A	434	ALA	2.5
1	B	440	ASN	2.5
1	B	437	VAL	2.4
1	B	46	LYS	2.4
1	A	56	PRO	2.4
1	A	263	LEU	2.3
1	B	68	PRO	2.3
1	A	394	TYR	2.3
1	A	208	GLU	2.3
1	A	234	VAL	2.3
1	A	241	PRO	2.3
1	A	52	PHE	2.3
1	B	505	TYR	2.3
1	A	441	SER	2.2
1	B	373	HIS	2.2
1	A	523	ARG	2.2
1	A	515	PRO	2.1
1	A	444	GLU	2.1
1	A	211	ARG	2.1
1	A	245	ARG	2.1
1	A	236	PRO	2.1
1	B	67	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	431	ALA	2.1
1	B	232	ASN	2.1
1	A	267	ARG	2.1
1	B	231	ILE	2.1
1	B	185	VAL	2.1
1	A	218	ALA	2.1
1	A	430	ALA	2.1
1	A	121	LEU	2.1
1	A	207	ASP	2.0
1	A	428	ASP	2.0
1	A	229	THR	2.0
1	A	528	PHE	2.0
1	B	442	SER	2.0
1	B	263	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

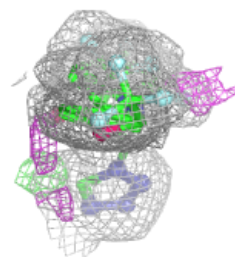
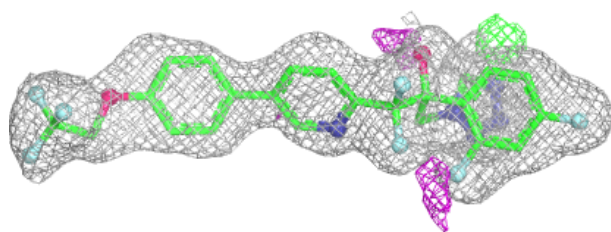
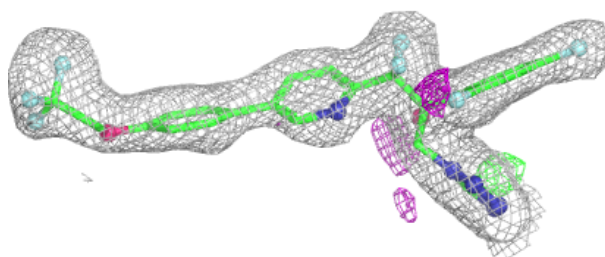
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	VT1	A	602	37/37	0.92	0.10	21,30,69,71	0
3	VT1	B	602	37/37	0.94	0.10	20,30,82,89	0
2	HEM	A	601	43/43	0.97	0.07	20,25,30,31	0
2	HEM	B	601	43/43	0.97	0.07	18,25,28,31	0

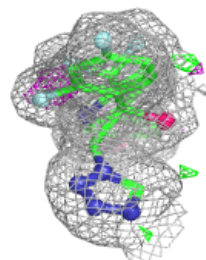
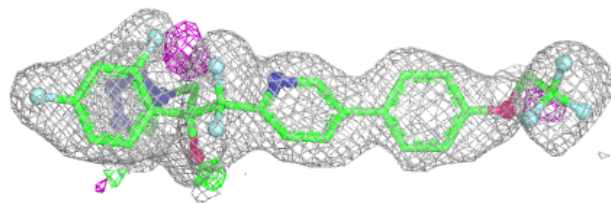
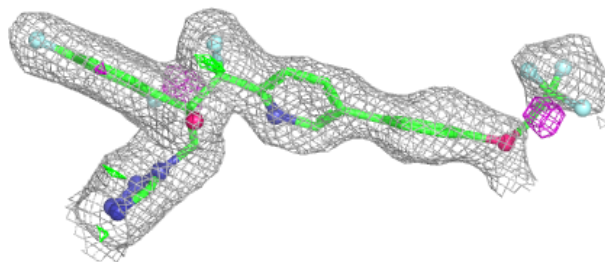
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around VT1 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)

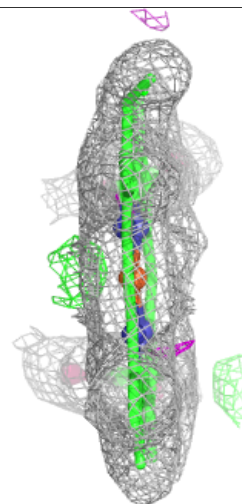
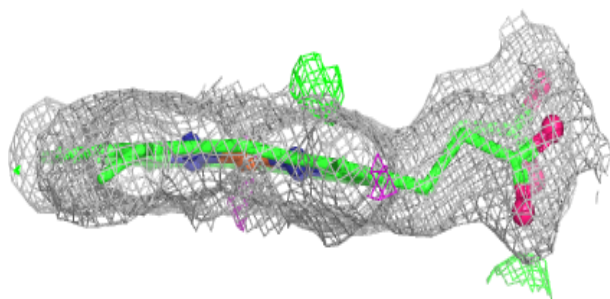
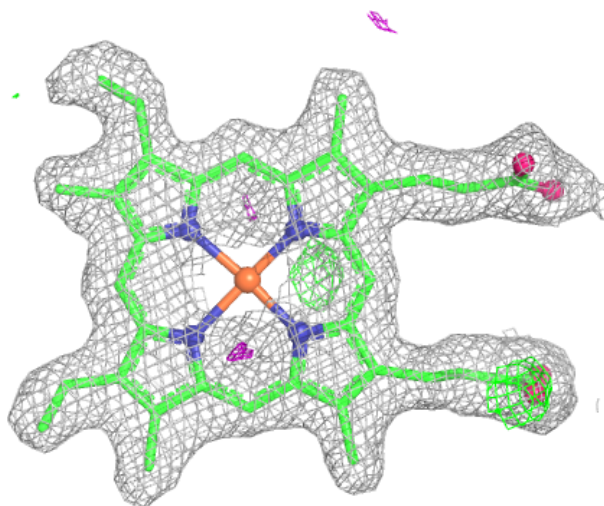
**Electron density around VT1 B 602:**

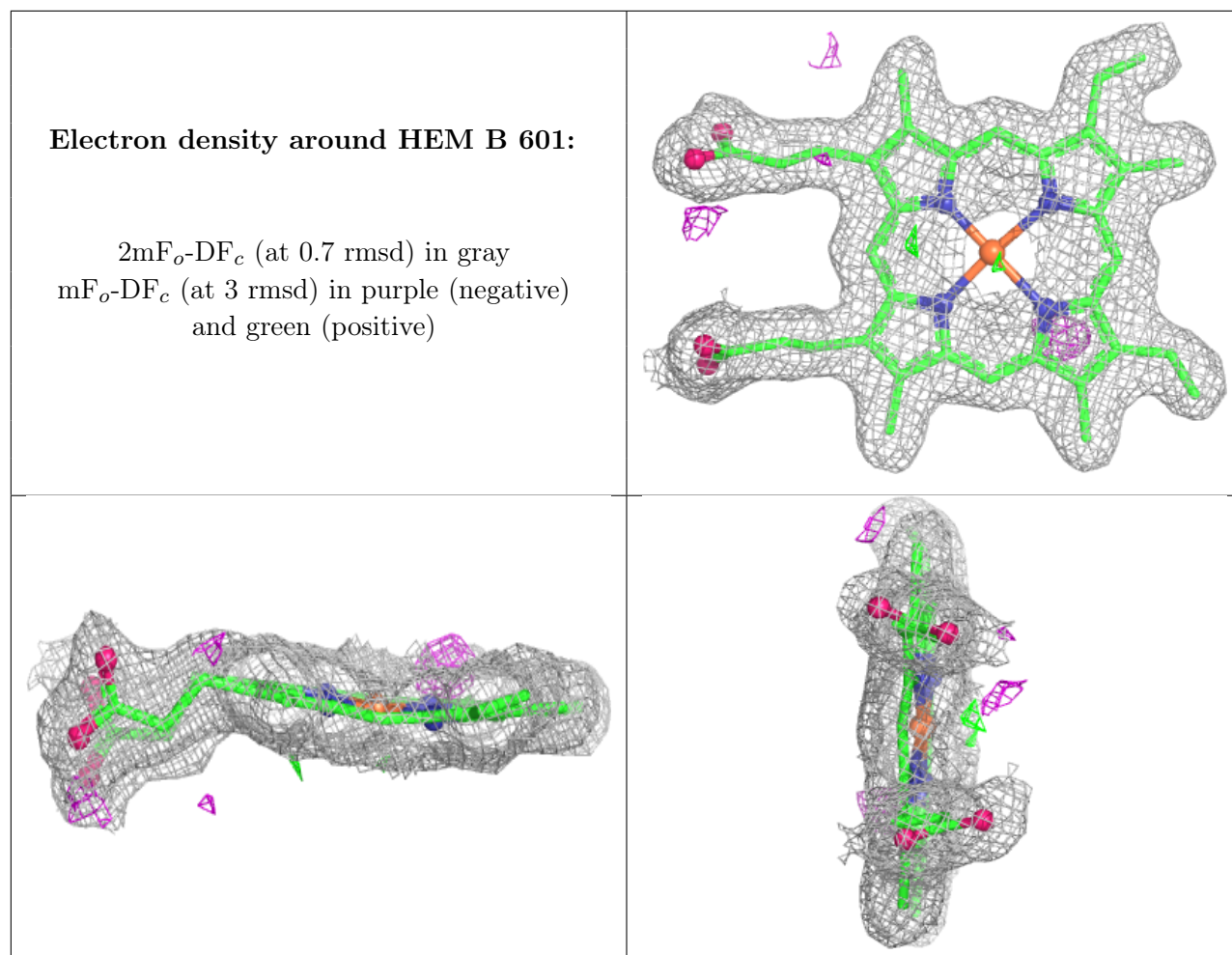
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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 ⓘ

There are no such residues in this entry.