



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:24 PM UTC

PDB ID : 1N6B / pdb_00001n6b
Title : Microsomal Cytochrome P450 2C5/3LVdH Complex with a dimethyl derivative of sulfaphenazole
Authors : Wester, M.R.; Johnson, E.F.; Marques-Soares, C.; Dansette, P.M.; Mansuy, D.; Stout, C.D.
Deposited on : 2002-11-09
Resolution : 2.30 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

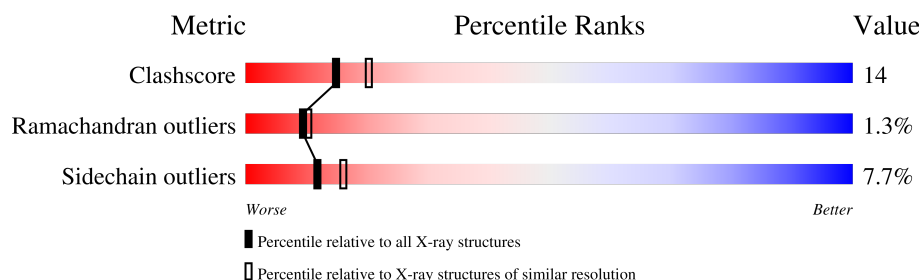
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	473	68% 25% . .

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	DMZ	A	501[B]	X	-	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3914 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 2C5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	462	3697	2384	619	672	22	0	0	0

There are 17 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	MET	-	engineered mutation	UNP P00179
A	20	ALA	-	engineered mutation	UNP P00179
A	21	LYS	-	engineered mutation	UNP P00179
A	22	LYS	-	engineered mutation	UNP P00179
A	23	THR	-	engineered mutation	UNP P00179
A	24	SER	-	engineered mutation	UNP P00179
A	25	SER	-	engineered mutation	UNP P00179
A	26	LYS	-	engineered mutation	UNP P00179
A	202	HIS	ASN	engineered mutation	UNP P00179
A	206	GLU	ARG	engineered mutation	UNP P00179
A	207	LEU	ILE	engineered mutation	UNP P00179
A	209	GLY	SER	engineered mutation	UNP P00179
A	210	THR	SER	engineered mutation	UNP P00179
A	488	HIS	-	expression tag	UNP P00179
A	489	HIS	-	expression tag	UNP P00179
A	490	HIS	-	expression tag	UNP P00179
A	491	HIS	-	expression tag	UNP P00179

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



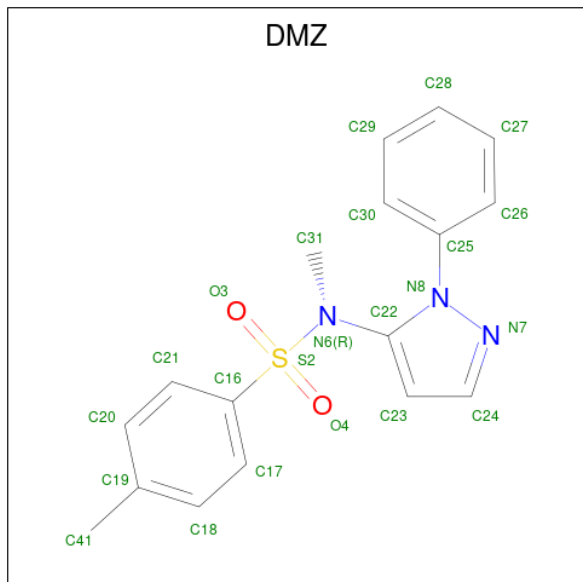
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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



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

- Molecule 4 is 4-METHYL-N-METHYL-N-(2-PHENYL-2H-PYRAZOL-3-YL)BENZENESULFONAMIDE (CCD ID: DMZ) (formula: $C_{17}H_{17}N_3O_2S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	1
			46	34	6	4	2		

- Molecule 5 is water.

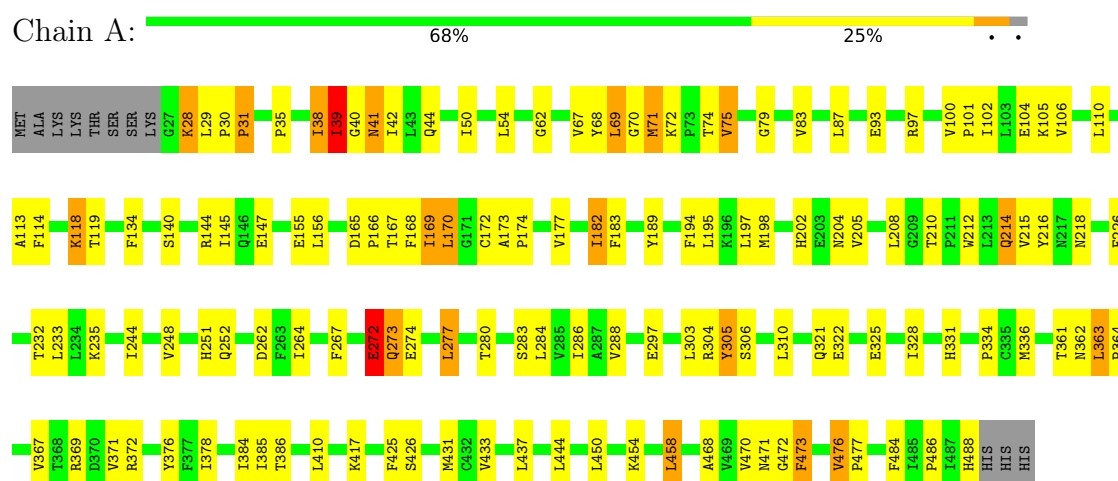
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	118	Total	O	0	0
			118	118		

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. 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. 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.

Note EDS was not executed.

• Molecule 1: Cytochrome P450 2C5



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	74.33Å 134.29Å 171.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30	Depositor
% Data completeness (in resolution range)	96.6 (50.00-2.30)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.257 , 0.292	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3914	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: DMZ, SO4, 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	0.57	1/3785 (0.0%)	1.06	23/5123 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	361	THR	CA-CB	5.00	1.60	1.53

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	363	LEU	N-CA-C	-7.86	99.16	110.40
1	A	69	LEU	N-CA-C	-7.28	104.93	113.88
1	A	262	ASP	N-CA-C	6.81	117.23	108.24
1	A	62	GLY	CA-C-N	6.79	126.58	119.05
1	A	62	GLY	C-N-CA	6.79	126.58	119.05
1	A	41	ASN	N-CA-C	6.74	120.81	112.59
1	A	454	LYS	N-CA-C	-6.71	98.46	109.40
1	A	476	VAL	CA-C-N	6.35	124.23	119.66
1	A	476	VAL	C-N-CA	6.35	124.23	119.66
1	A	100	VAL	CA-C-N	5.87	125.34	119.24
1	A	100	VAL	C-N-CA	5.87	125.34	119.24
1	A	40	GLY	N-CA-C	5.85	127.04	113.18
1	A	39	ILE	N-CA-C	5.82	121.44	109.34
1	A	331	HIS	N-CA-C	5.71	120.01	111.96
1	A	113	ALA	N-CA-C	5.68	117.55	111.36
1	A	168	PHE	N-CA-C	5.54	117.00	111.07
1	A	182	ILE	N-CA-C	5.47	116.54	111.45
1	A	165	ASP	CA-C-N	5.38	125.46	119.87
1	A	165	ASP	C-N-CA	5.38	125.46	119.87
1	A	336	MET	N-CA-C	5.31	118.55	111.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	226	PHE	CA-C-N	5.09	124.87	119.28
1	A	226	PHE	C-N-CA	5.09	124.87	119.28
1	A	106	VAL	N-CA-C	5.05	115.26	110.42

There are no chirality outliers.

There are no planarity outliers.

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	3697	0	3724	105	0
2	A	10	0	0	0	0
3	A	43	0	30	2	0
4	A	46	0	34	0	0
5	A	118	0	0	6	0
All	All	3914	0	3788	105	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 14.

All (105) 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:182:ILE:HD13	1:A:288:VAL:HG13	1.47	0.94
1:A:41:ASN:HD22	1:A:44:GLN:NE2	1.71	0.89
1:A:41:ASN:HD22	1:A:44:GLN:HE22	1.20	0.88
1:A:172:CYS:SG	1:A:198:MET:HE1	2.18	0.84
1:A:101:PRO:HG2	1:A:102:ILE:HD12	1.58	0.82
1:A:277:LEU:H	1:A:277:LEU:HD23	1.46	0.80
1:A:38:ILE:HG22	1:A:39:ILE:HG23	1.68	0.75
1:A:214:GLN:HE21	1:A:214:GLN:H	1.33	0.74
1:A:205:VAL:HG11	1:A:297:GLU:HG2	1.70	0.73
1:A:102:ILE:HD12	1:A:102:ILE:H	1.52	0.73
1:A:101:PRO:CG	5:A:819:HOH:O	2.37	0.72
1:A:30:PRO:HA	1:A:376:TYR:CE2	2.23	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:LYS:HE3	5:A:817:HOH:O	1.91	0.71
1:A:473:PHE:CD1	1:A:473:PHE:N	2.57	0.70
1:A:105:LYS:CE	5:A:817:HOH:O	2.41	0.69
1:A:101:PRO:HG2	5:A:819:HOH:O	1.95	0.67
1:A:69:LEU:HB2	1:A:218:ASN:HD22	1.59	0.66
1:A:166:PRO:HB2	1:A:170:LEU:HD22	1.80	0.63
1:A:41:ASN:HA	1:A:44:GLN:HE21	1.63	0.63
1:A:134:PHE:CE1	1:A:140:SER:HB3	2.34	0.62
1:A:42:ILE:H	1:A:42:ILE:HD12	1.65	0.61
1:A:156:LEU:CD2	1:A:169:ILE:HD11	2.30	0.60
1:A:71:MET:H	1:A:71:MET:HE3	1.67	0.59
1:A:328:ILE:HD13	1:A:334:PRO:HB3	1.85	0.58
1:A:484:PHE:O	1:A:486:PRO:HD3	2.04	0.58
1:A:473:PHE:N	1:A:473:PHE:HD1	2.02	0.57
1:A:322:GLU:O	1:A:325:GLU:HG3	2.04	0.57
1:A:212:TRP:CE2	1:A:385:ILE:HD13	2.39	0.56
1:A:42:ILE:HD12	1:A:42:ILE:N	2.21	0.56
1:A:69:LEU:HD22	1:A:69:LEU:N	2.21	0.55
1:A:214:GLN:H	1:A:214:GLN:NE2	2.03	0.55
1:A:39:ILE:O	1:A:42:ILE:HD13	2.06	0.55
1:A:155:GLU:HG3	1:A:189:TYR:CE1	2.41	0.55
1:A:31:PRO:HD3	1:A:376:TYR:CE2	2.42	0.55
1:A:118:LYS:HG2	1:A:119:THR:H	1.71	0.54
1:A:177:VAL:HG11	1:A:437:LEU:HD11	1.89	0.54
1:A:310:LEU:HD13	1:A:450:LEU:HD11	1.89	0.54
1:A:458:LEU:HD12	1:A:458:LEU:N	2.23	0.54
1:A:362:ASN:ND2	1:A:363:LEU:H	2.06	0.54
1:A:38:ILE:HG22	1:A:39:ILE:N	2.22	0.54
1:A:321:GLN:HG2	1:A:450:LEU:O	2.08	0.53
1:A:38:ILE:HA	1:A:71:MET:HE2	1.89	0.53
1:A:102:ILE:HD12	1:A:102:ILE:N	2.23	0.53
1:A:437:LEU:HD12	1:A:437:LEU:O	2.09	0.53
1:A:232:THR:HA	1:A:235:LYS:HE3	1.90	0.53
1:A:173:ALA:HB3	1:A:174:PRO:HD3	1.91	0.52
1:A:458:LEU:H	1:A:458:LEU:CD1	2.23	0.52
1:A:172:CYS:HA	1:A:198:MET:SD	2.49	0.52
1:A:42:ILE:H	1:A:42:ILE:CD1	2.22	0.52
1:A:362:ASN:HD22	1:A:363:LEU:H	1.58	0.52
1:A:202:HIS:O	1:A:205:VAL:HG12	2.09	0.51
1:A:167:THR:HA	1:A:304:ARG:HD3	1.91	0.51
1:A:97:ARG:HG2	1:A:114:PHE:HA	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:LYS:HG2	1:A:119:THR:N	2.26	0.51
1:A:177:VAL:CG1	1:A:437:LEU:HD11	2.41	0.50
1:A:28:LYS:HG3	1:A:29:LEU:N	2.27	0.49
1:A:87:LEU:HD11	1:A:386:THR:HG21	1.93	0.49
1:A:155:GLU:HG3	1:A:189:TYR:CD1	2.47	0.49
1:A:458:LEU:HD12	1:A:458:LEU:H	1.79	0.48
1:A:305:TYR:OH	1:A:468:ALA:HB2	2.13	0.48
1:A:118:LYS:N	1:A:118:LYS:HD2	2.29	0.48
1:A:97:ARG:HD3	1:A:364:PRO:O	2.14	0.47
1:A:267:PHE:CE2	1:A:284:LEU:HB2	2.49	0.47
1:A:93:GLU:HB3	1:A:371:VAL:HG11	1.96	0.47
1:A:41:ASN:ND2	1:A:44:GLN:HE22	2.00	0.47
1:A:54:LEU:HD11	1:A:212:TRP:HZ2	1.80	0.47
1:A:210:THR:N	1:A:472:GLY:HA2	2.29	0.47
1:A:74:THR:HG21	1:A:385:ILE:HD12	1.97	0.46
1:A:28:LYS:HE2	1:A:29:LEU:H	1.81	0.46
1:A:145:ILE:CG2	1:A:444:LEU:HD12	2.45	0.46
1:A:280:THR:H	1:A:283:SER:HB3	1.80	0.46
1:A:426:SER:CB	3:A:500:HEM:HBA1	2.45	0.46
1:A:118:LYS:HD2	1:A:118:LYS:H	1.80	0.46
1:A:215:VAL:HG13	1:A:216:TYR:N	2.31	0.46
1:A:362:ASN:HD22	1:A:363:LEU:N	2.14	0.46
1:A:371:VAL:HG23	1:A:378:ILE:HB	1.96	0.45
1:A:104:GLU:HG2	5:A:822:HOH:O	2.17	0.45
1:A:244:ILE:O	1:A:248:VAL:HG23	2.16	0.45
1:A:41:ASN:ND2	1:A:44:GLN:NE2	2.53	0.45
1:A:297:GLU:HB2	5:A:613:HOH:O	2.16	0.45
1:A:79:GLY:O	1:A:83:VAL:HG23	2.17	0.44
1:A:426:SER:HB3	3:A:500:HEM:HBA1	2.00	0.44
1:A:156:LEU:HD22	1:A:484:PHE:CD1	2.53	0.44
1:A:183:PHE:CD1	1:A:264:ILE:HD11	2.53	0.44
1:A:272:GLU:HB3	1:A:273:GLN:H	1.59	0.44
1:A:425:PHE:O	1:A:426:SER:HB3	2.17	0.44
1:A:248:VAL:O	1:A:252:GLN:HG3	2.16	0.43
1:A:102:ILE:HG21	1:A:233:LEU:HD13	2.00	0.43
1:A:367:VAL:HG21	1:A:378:ILE:HG22	2.01	0.43
1:A:194:PHE:CD2	1:A:198:MET:HE2	2.54	0.43
1:A:194:PHE:CE2	1:A:198:MET:HE2	2.53	0.42
1:A:277:LEU:HD23	1:A:277:LEU:N	2.26	0.42
1:A:232:THR:HG23	1:A:235:LYS:NZ	2.33	0.42
1:A:458:LEU:N	1:A:458:LEU:CD1	2.81	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:ILE:H	1:A:102:ILE:CD1	2.25	0.42
1:A:110:LEU:HD12	1:A:286:ILE:HD12	2.01	0.42
1:A:35:PRO:HB3	1:A:68:TYR:CD1	2.56	0.41
1:A:70:GLY:C	1:A:72:LYS:H	2.29	0.41
1:A:214:GLN:OE1	1:A:385:ILE:HD11	2.21	0.41
1:A:183:PHE:HD1	1:A:264:ILE:HD11	1.85	0.41
1:A:75:VAL:CG1	1:A:384:ILE:HD13	2.50	0.41
1:A:30:PRO:HA	1:A:31:PRO:HD3	1.91	0.41
1:A:101:PRO:HG2	1:A:102:ILE:CD1	2.41	0.41
1:A:67:VAL:HG23	1:A:74:THR:HB	2.03	0.41
1:A:303:LEU:O	1:A:306:SER:HB2	2.21	0.41

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	460/473 (97%)	430 (94%)	24 (5%)	6 (1%)	9 10

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	38	ILE
1	A	273	GLN
1	A	31	PRO
1	A	272	GLU
1	A	471	ASN
1	A	39	ILE

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	415/425 (98%)	383 (92%)	32 (8%)	12	16

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LYS
1	A	39	ILE
1	A	50	ILE
1	A	71	MET
1	A	75	VAL
1	A	118	LYS
1	A	144	ARG
1	A	147	GLU
1	A	169	ILE
1	A	170	LEU
1	A	195	LEU
1	A	197	LEU
1	A	204	ASN
1	A	208	LEU
1	A	214	GLN
1	A	251	HIS
1	A	272	GLU
1	A	274	GLU
1	A	277	LEU
1	A	305	TYR
1	A	369	ARG
1	A	372	ARG
1	A	410	LEU
1	A	417	LYS
1	A	431	MET
1	A	433	VAL
1	A	458	LEU
1	A	470	VAL
1	A	473	PHE
1	A	476	VAL

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Mol	Chain	Res	Type
1	A	477	PRO
1	A	488	HIS

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	44	GLN
1	A	184	HIS
1	A	185	ASN
1	A	214	GLN
1	A	218	ASN
1	A	273	GLN
1	A	313	HIS
1	A	321	GLN
1	A	362	ASN
1	A	471	ASN

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](#)

5 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	HEM	A	500	1	50,50,50	2.45	27 (54%)	67,82,82	1.29	10 (14%)
4	DMZ	A	501[A]	-	24,25,25	3.49	9 (37%)	32,36,36	1.36	4 (12%)
2	SO4	A	503	-	4,4,4	0.39	0	6,6,6	0.16	0
4	DMZ	A	501[B]	-	24,25,25	3.48	7 (29%)	32,36,36	1.34	4 (12%)
2	SO4	A	502	-	4,4,4	0.37	0	6,6,6	0.22	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HEM	A	500	1	-	2/14/54/54	-
4	DMZ	A	501[B]	-	1/1/2/2	3/20/20/20	0/3/3/3
4	DMZ	A	501[A]	-	-	2/20/20/20	0/3/3/3

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	501[B]	DMZ	O3-S2	9.04	1.53	1.43
4	A	501[B]	DMZ	O4-S2	8.98	1.53	1.43
4	A	501[A]	DMZ	O4-S2	8.88	1.53	1.43
4	A	501[A]	DMZ	O3-S2	8.81	1.52	1.43
4	A	501[A]	DMZ	C22-N8	7.70	1.43	1.36
4	A	501[B]	DMZ	C22-N8	7.17	1.42	1.36
3	A	500	HEM	C1D-ND	-4.78	1.29	1.38
4	A	501[A]	DMZ	S2-N6	4.69	1.76	1.65
3	A	500	HEM	C1B-NB	-4.55	1.32	1.40
4	A	501[B]	DMZ	S2-N6	4.34	1.75	1.65
3	A	500	HEM	C4A-NA	-4.28	1.31	1.39
3	A	500	HEM	C4D-ND	-4.14	1.33	1.40
3	A	500	HEM	C1C-NC	-4.07	1.31	1.39
3	A	500	HEM	CAC-C3C	-4.07	1.36	1.47
3	A	500	HEM	C4B-NB	-3.94	1.31	1.38
3	A	500	HEM	C1C-C2C	-3.88	1.37	1.45
3	A	500	HEM	CBB-CAB	3.68	1.48	1.30
4	A	501[B]	DMZ	N8-N7	-3.35	1.33	1.38
3	A	500	HEM	C4C-NC	-3.10	1.33	1.39
3	A	500	HEM	C3C-C4C	-2.92	1.40	1.46
3	A	500	HEM	C4D-C3D	-2.89	1.40	1.45
3	A	500	HEM	CHD-C1D	-2.89	1.32	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	HEM	FE-ND	2.88	2.03	1.94
3	A	500	HEM	O2A-CGA	-2.75	1.21	1.30
3	A	500	HEM	C1A-C2A	-2.75	1.39	1.44
4	A	501[A]	DMZ	C23-C24	2.67	1.44	1.39
3	A	500	HEM	C1D-C2D	-2.59	1.39	1.44
3	A	500	HEM	CMA-C3A	2.53	1.56	1.50
3	A	500	HEM	FE-NB	2.51	2.02	1.94
4	A	501[A]	DMZ	N8-N7	-2.50	1.34	1.38
3	A	500	HEM	C3B-C2B	2.49	1.42	1.37
3	A	500	HEM	FE-NA	2.49	2.03	1.95
3	A	500	HEM	CBD-CGD	2.39	1.56	1.50
3	A	500	HEM	CBC-CAC	2.34	1.41	1.30
3	A	500	HEM	C3C-C2C	2.31	1.41	1.37
4	A	501[B]	DMZ	C17-C16	2.28	1.42	1.38
3	A	500	HEM	C3B-C4B	-2.25	1.40	1.44
4	A	501[B]	DMZ	C23-C24	2.18	1.43	1.39
3	A	500	HEM	CMC-C2C	2.11	1.55	1.50
4	A	501[A]	DMZ	C17-C16	2.04	1.42	1.38
4	A	501[A]	DMZ	C21-C16	2.03	1.42	1.38
3	A	500	HEM	FE-NC	2.02	2.01	1.95
4	A	501[A]	DMZ	C26-C25	2.01	1.43	1.39

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	501[A]	DMZ	O4-S2-O3	-4.42	112.69	119.59
4	A	501[B]	DMZ	O4-S2-O3	-4.42	112.70	119.59
4	A	501[B]	DMZ	C24-N7-N8	3.33	107.57	104.25
4	A	501[A]	DMZ	C24-N7-N8	3.05	107.29	104.25
4	A	501[A]	DMZ	C23-C22-N8	-2.93	105.62	107.14
3	A	500	HEM	C4C-C3C-C2C	-2.91	104.29	106.81
4	A	501[B]	DMZ	C23-C22-N8	-2.79	105.69	107.14
3	A	500	HEM	C1B-NB-C4B	2.72	108.43	105.21
3	A	500	HEM	C3B-C4B-NB	2.68	111.39	109.47
4	A	501[A]	DMZ	O4-S2-C16	2.67	111.43	108.10
3	A	500	HEM	C4B-C3B-C2B	-2.62	104.87	107.28
3	A	500	HEM	C3B-C2B-C1B	-2.55	104.50	106.41
3	A	500	HEM	CHD-C4C-NC	-2.43	121.80	124.45
3	A	500	HEM	O2A-CGA-CBA	2.34	121.39	114.00
4	A	501[B]	DMZ	O4-S2-C16	2.19	110.83	108.10
3	A	500	HEM	CHC-C1C-NC	-2.09	122.17	124.45
3	A	500	HEM	C1C-CHC-C4B	2.08	130.45	126.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	500	HEM	C4D-ND-C1D	2.08	107.67	105.21

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	A	501[B]	DMZ	N6

All (7) torsion outliers are listed below:

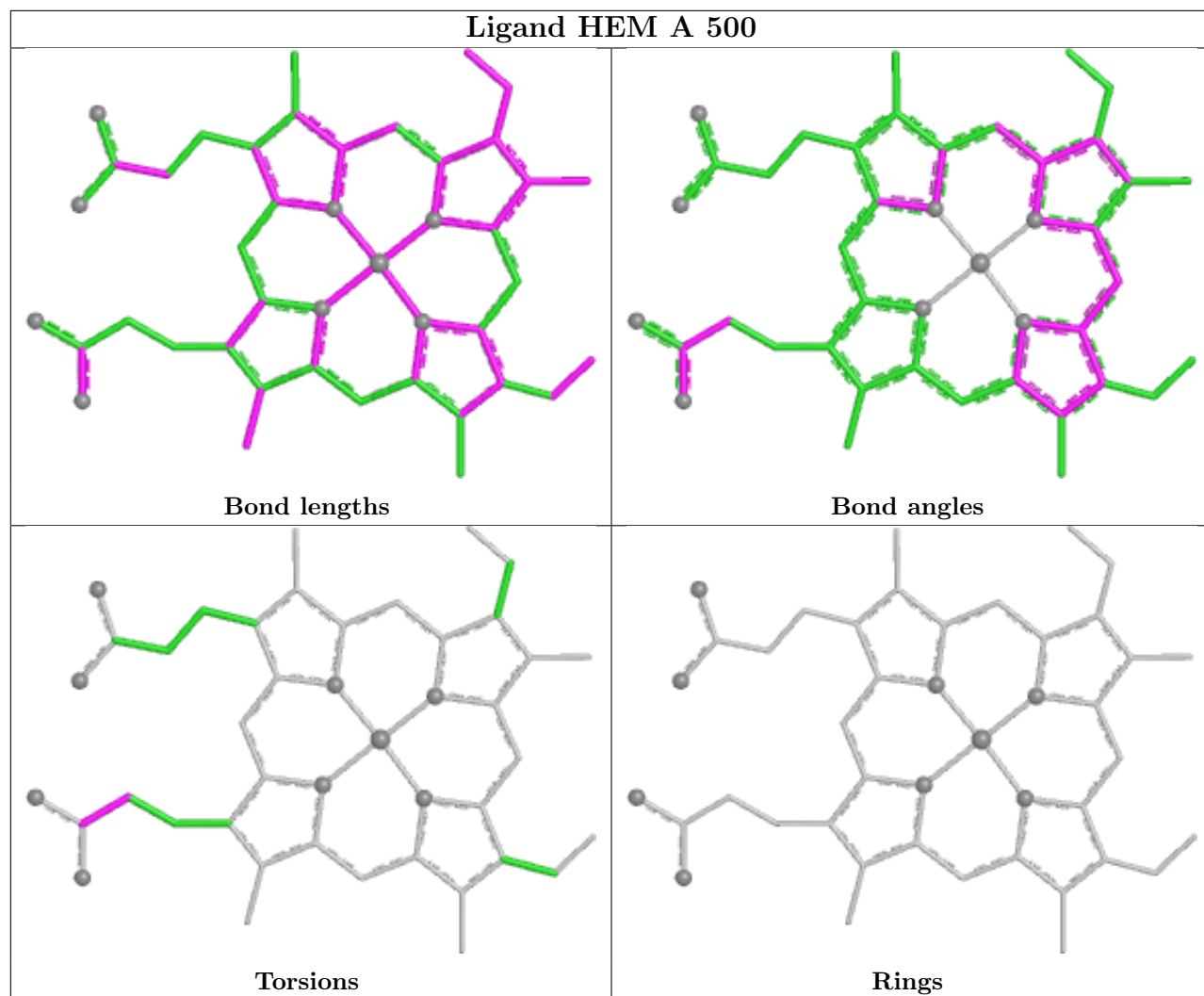
Mol	Chain	Res	Type	Atoms
4	A	501[A]	DMZ	C23-C22-N6-S2
4	A	501[A]	DMZ	N8-C22-N6-S2
4	A	501[B]	DMZ	C22-N6-S2-C16
4	A	501[B]	DMZ	C22-N6-S2-O3
3	A	500	HEM	CAA-CBA-CGA-O1A
3	A	500	HEM	CAA-CBA-CGA-O2A
4	A	501[B]	DMZ	C31-N6-S2-O3

There are no ring outliers.

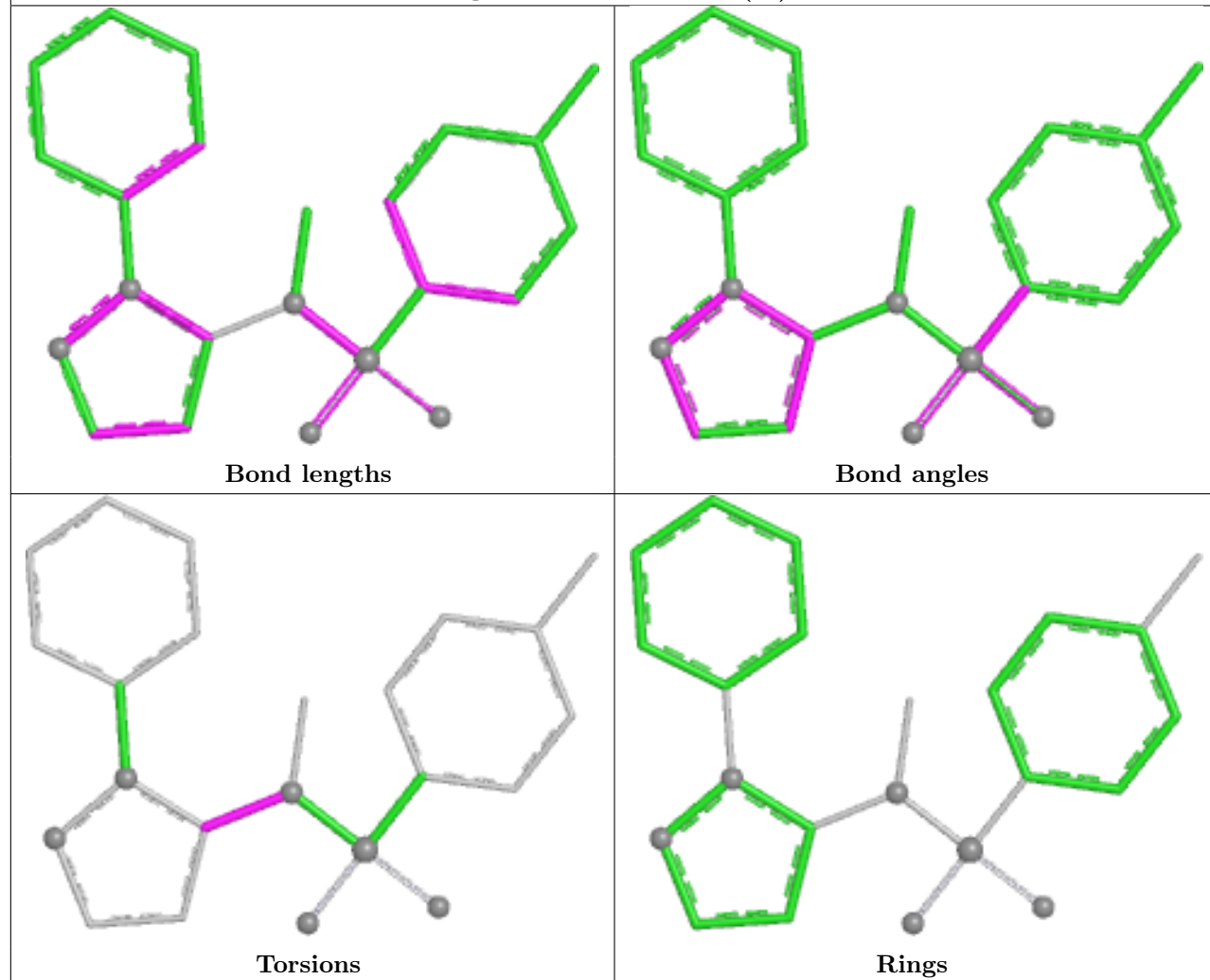
1 monomer is involved in 2 short contacts:

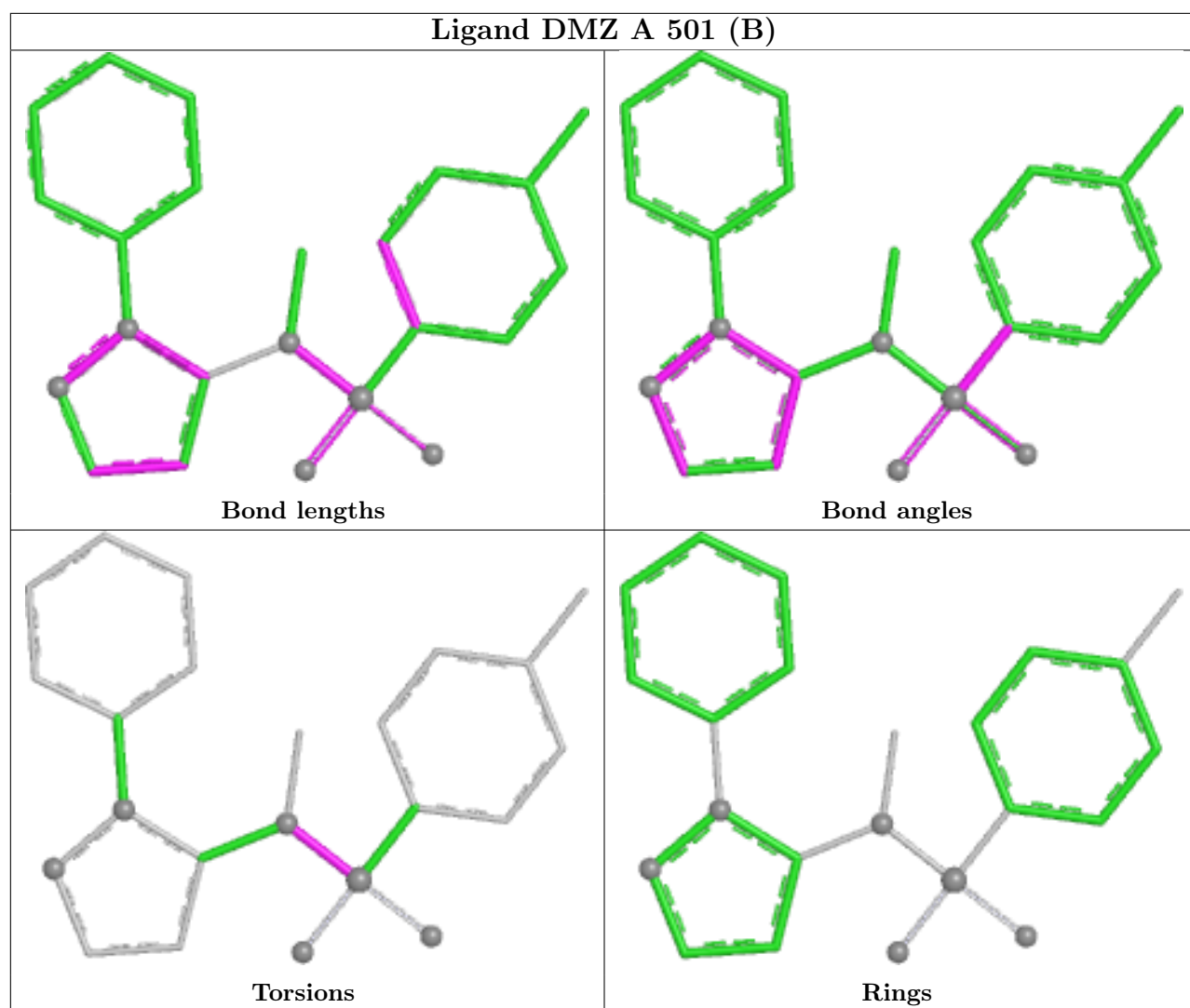
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	500	HEM	2	0

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 DMZ A 501 (A)





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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.