



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 01:59 AM UTC

PDB ID : 2PO5 / pdb_00002po5
Title : Crystal structure of human ferrochelatase mutant with His 263 replaced by Cys
Authors : Dailey, H.A.; Wu, C.-K.; Horanyi, P.; Medlock, A.E.; Najahi-Missaoui, A.E.W.; Burden, A.; Dailey, T.A.; Rose, J.P.
Deposited on : 2007-04-25
Resolution : 2.20 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

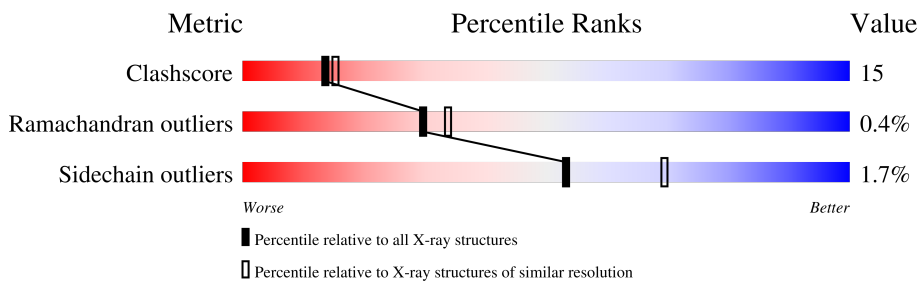
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)

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	359	 70% 28% .
1	B	359	 75% 24% .

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6192 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.

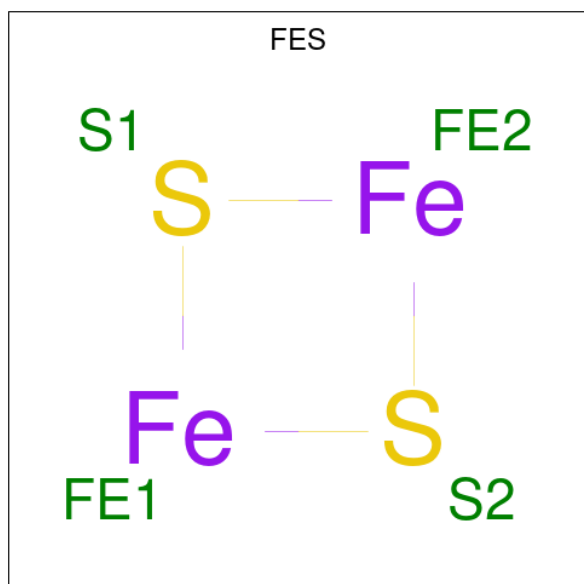
- Molecule 1 is a protein called Ferrochelatase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	359	Total	C	N	O	S	0	0	0
			2851	1817	488	527	19			
1	B	359	Total	C	N	O	S	0	0	0
			2851	1817	488	527	19			

There are 4 discrepancies between the modelled and reference sequences:

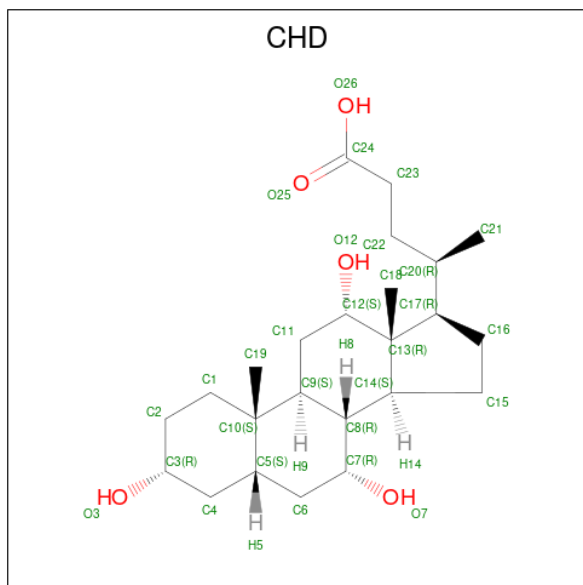
Chain	Residue	Modelled	Actual	Comment	Reference
A	115	LEU	ARG	engineered mutation	UNP P22830
A	263	CYS	HIS	engineered mutation	UNP P22830
B	115	LEU	ARG	engineered mutation	UNP P22830
B	263	CYS	HIS	engineered mutation	UNP P22830

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Fe S 4 2 2	0	0
2	B	1	Total Fe S 4 2 2	0	0

- Molecule 3 is CHOLIC ACID (CCD ID: CHD) (formula: $C_{24}H_{40}O_5$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 29 24 5	0	0
3	A	1	Total C O 29 24 5	0	0
3	A	1	Total C O 29 24 5	0	0
3	B	1	Total C O 29 24 5	0	0
3	B	1	Total C O 29 24 5	0	0
3	B	1	Total C O 29 24 5	0	0

- Molecule 4 is water.

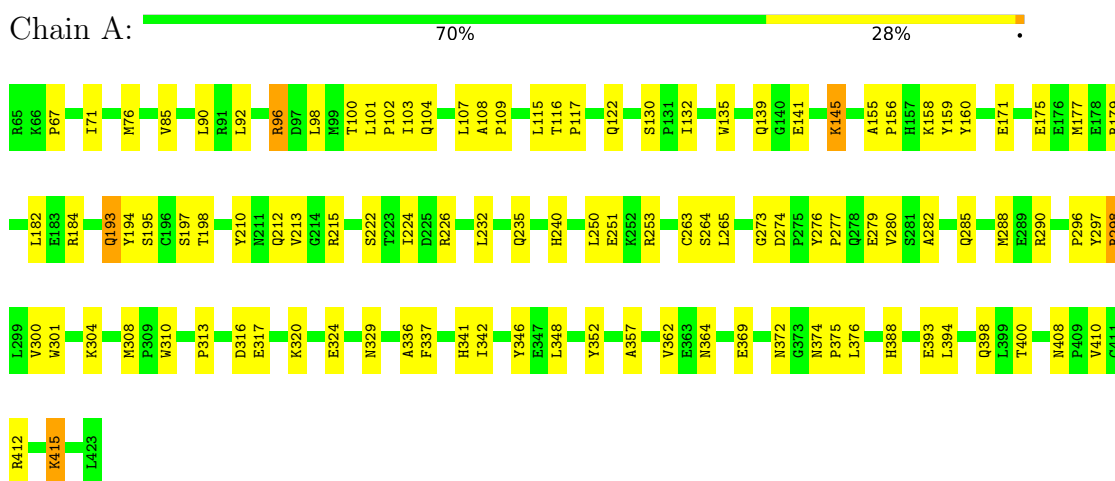
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	157	Total O 157 157	0	0
4	B	151	Total O 151 151	0	0

3 Residue-property plots

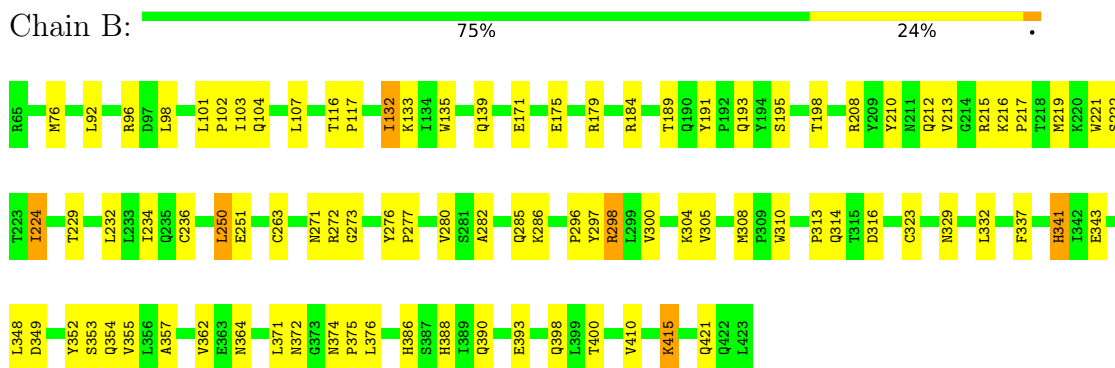
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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



- Molecule 1: Ferrochelatase, mitochondrial



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	93.42Å 87.73Å 109.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.90 – 2.20	Depositor
% Data completeness (in resolution range)	96.3 (19.90-2.20)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.209 , 0.240	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6192	wwPDB-VP
Average B, all atoms (Å ²)	26.0	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: CHD, FES

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	0.91	1/2920 (0.0%)	0.86	6/3965 (0.2%)
1	B	0.43	1/2920 (0.0%)	0.84	2/3965 (0.1%)
All	All	0.71	2/5840 (0.0%)	0.85	8/7930 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	96	ARG	CD-NE	44.14	2.08	1.46
1	B	96	ARG	CD-NE	-9.27	1.33	1.46

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	274	ASP	CA-C-N	7.13	127.12	119.28
1	A	274	ASP	C-N-CA	7.13	127.12	119.28
1	B	224	ILE	N-CA-C	-6.49	95.84	109.34
1	A	224	ILE	N-CA-C	-6.22	96.41	109.34
1	B	132	ILE	N-CA-C	5.83	117.27	110.62
1	A	194	TYR	N-CA-C	5.57	118.32	109.96
1	A	85	VAL	N-CA-C	5.49	115.69	110.42
1	A	394	LEU	N-CA-C	-5.20	107.06	113.41

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	2851	0	2820	92	0
1	B	2851	0	2820	79	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
3	A	87	0	117	13	0
3	B	87	0	117	10	0
4	A	157	0	0	2	0
4	B	151	0	0	1	0
All	All	6192	0	5874	172	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 15.

All (172) 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:96:ARG:CD	1:A:96:ARG:NE	2.08	1.16
1:A:297:TYR:H	1:B:398:GLN:HE22	1.13	0.97
1:B:250:LEU:HD12	1:B:250:LEU:H	1.30	0.94
1:B:337:PHE:CE2	1:B:341:HIS:HE1	1.86	0.92
1:A:398:GLN:HE22	1:B:297:TYR:H	1.19	0.90
1:B:337:PHE:CE1	1:B:341:HIS:CE1	2.59	0.90
1:B:337:PHE:CE1	1:B:341:HIS:NE2	2.41	0.89
1:B:229:THR:HB	1:B:286:LYS:HE2	1.57	0.86
1:B:337:PHE:CZ	1:B:341:HIS:CE1	2.64	0.85
1:A:250:LEU:HD12	1:A:250:LEU:H	1.42	0.83
1:A:213:VAL:HG13	1:A:215:ARG:HG2	1.60	0.82
1:B:337:PHE:CD1	1:B:341:HIS:CE1	2.67	0.82
1:B:232:LEU:HD22	1:B:376:LEU:HD11	1.61	0.81
1:B:101:LEU:HB2	1:B:104:GLN:HG3	1.63	0.80
1:B:337:PHE:CD2	1:B:341:HIS:HE1	1.99	0.80
1:B:337:PHE:CZ	1:B:341:HIS:HE1	2.00	0.79
1:B:323:CYS:SG	1:B:362:VAL:HG12	2.24	0.78
1:B:400:THR:HA	1:B:415:LYS:HD2	1.66	0.77
1:A:98:LEU:HD11	3:A:501:CHD:H162	1.67	0.76
1:B:132:ILE:HD12	1:B:133:LYS:N	1.99	0.76
1:A:193:GLN:HG3	1:A:280:VAL:HG22	1.67	0.76
1:B:337:PHE:CD1	1:B:341:HIS:NE2	2.54	0.76
1:B:308:MET:HG3	3:B:502:CHD:O3	1.87	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:MET:HE1	3:A:501:CHD:H151	1.70	0.74
1:A:265:LEU:HD22	3:A:501:CHD:O25	1.88	0.73
1:A:341:HIS:CD2	1:A:342:ILE:H	2.06	0.73
1:B:221:TRP:O	1:B:421:GLN:HG2	1.89	0.72
1:A:222:SER:OG	1:A:388:HIS:HE1	1.73	0.71
1:B:337:PHE:CG	1:B:341:HIS:CE1	2.80	0.69
1:A:357:ALA:HB1	1:A:362:VAL:HG11	1.76	0.67
1:B:374:ASN:ND2	1:B:376:LEU:H	1.93	0.66
1:B:250:LEU:H	1:B:250:LEU:CD1	2.06	0.66
1:A:197:SER:HB3	3:A:501:CHD:O26	1.95	0.66
1:A:297:TYR:N	1:B:398:GLN:HE22	1.89	0.66
1:B:341:HIS:NE2	1:B:343:GLU:HB2	2.11	0.65
1:B:213:VAL:HG21	1:B:215:ARG:HH21	1.61	0.65
1:A:130:SER:OG	1:A:132:ILE:HG13	1.98	0.64
1:B:263:CYS:HG	1:B:337:PHE:HZ	1.43	0.64
1:A:308:MET:HG3	3:A:502:CHD:O3	1.98	0.64
1:B:222:SER:OG	1:B:388:HIS:HE1	1.80	0.64
1:B:374:ASN:HD22	1:B:374:ASN:C	2.05	0.63
1:B:337:PHE:CD2	1:B:341:HIS:CE1	2.84	0.63
1:A:240:HIS:HD2	1:A:369:GLU:O	1.82	0.62
1:A:132:ILE:HD11	1:A:341:HIS:CG	2.34	0.62
1:A:398:GLN:HE22	1:B:297:TYR:N	1.93	0.62
1:A:115:LEU:HD21	3:A:502:CHD:H231	1.81	0.62
1:A:341:HIS:CG	1:A:342:ILE:H	2.17	0.62
1:A:374:ASN:C	1:A:374:ASN:HD22	2.07	0.61
1:A:400:THR:HA	1:A:415:LYS:HD2	1.82	0.61
1:A:276:TYR:HB3	1:A:277:PRO:HD3	1.81	0.61
1:A:193:GLN:HG3	1:A:280:VAL:CG2	2.31	0.60
1:B:374:ASN:HD22	1:B:375:PRO:N	2.00	0.60
1:A:102:PRO:O	1:A:107:LEU:HD12	2.02	0.60
3:B:501:CHD:H211	3:B:502:CHD:H193	1.82	0.60
1:A:408:ASN:OD1	1:A:410:VAL:HG12	2.01	0.59
1:A:184:ARG:HH22	1:A:222:SER:CB	2.15	0.59
3:A:501:CHD:H212	3:A:501:CHD:H12	1.84	0.59
1:A:374:ASN:ND2	1:A:376:LEU:H	2.00	0.58
1:A:320:LYS:O	1:A:324:GLU:HG3	2.03	0.58
1:B:103:ILE:HG13	1:B:107:LEU:HG	1.85	0.58
1:A:374:ASN:HD22	1:A:375:PRO:N	2.00	0.58
3:B:501:CHD:C21	3:B:502:CHD:H193	2.34	0.58
1:B:76:MET:HB2	1:B:191:TYR:OH	2.04	0.57
1:B:316:ASP:HB3	1:B:352:TYR:CE1	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:501:CHD:H212	3:B:501:CHD:H12	1.86	0.57
1:A:184:ARG:HG3	4:A:642:HOH:O	2.04	0.57
1:A:329:ASN:HD22	1:A:364:ASN:HB2	1.70	0.57
1:B:234:ILE:HG13	1:B:286:LYS:HE3	1.86	0.57
1:B:337:PHE:CE2	1:B:341:HIS:CE1	2.77	0.57
1:A:300:VAL:HG12	1:A:313:PRO:HG2	1.86	0.56
1:A:132:ILE:HD11	1:A:341:HIS:CD2	2.40	0.56
1:A:285:GLN:HG2	1:B:282:ALA:HB1	1.87	0.56
1:B:98:LEU:HD11	3:B:501:CHD:H162	1.88	0.55
1:B:250:LEU:HD12	1:B:250:LEU:N	2.12	0.54
1:A:398:GLN:NE2	1:B:296:PRO:HA	2.22	0.54
1:B:276:TYR:HB3	1:B:277:PRO:HD3	1.90	0.54
1:B:357:ALA:HB1	1:B:362:VAL:HG21	1.89	0.54
3:B:501:CHD:H211	3:B:502:CHD:C19	2.38	0.54
1:A:304:LYS:HG3	1:A:310:TRP:CG	2.42	0.54
1:A:250:LEU:H	1:A:250:LEU:CD1	2.17	0.53
1:A:96:ARG:NE	1:A:104:GLN:NE2	2.56	0.53
1:A:235:GLN:HG3	1:A:290:ARG:CZ	2.39	0.53
1:A:316:ASP:HB3	1:A:352:TYR:CE1	2.44	0.53
1:B:341:HIS:CD2	1:B:343:GLU:HB2	2.44	0.53
1:A:116:THR:HB	1:A:117:PRO:HD3	1.90	0.52
1:B:236:CYS:HB3	1:B:371:LEU:HD22	1.90	0.52
1:A:282:ALA:HB1	1:B:285:GLN:HG2	1.90	0.52
1:A:415:LYS:NZ	1:A:415:LYS:HB3	2.24	0.52
1:A:296:PRO:HA	1:B:398:GLN:NE2	2.25	0.52
1:A:155:ALA:HB1	1:A:156:PRO:HA	1.92	0.52
1:B:102:PRO:O	1:B:103:ILE:C	2.53	0.52
1:B:263:CYS:SG	1:B:337:PHE:HZ	2.33	0.51
1:B:300:VAL:HG12	1:B:313:PRO:HG2	1.93	0.51
1:B:175:GLU:O	1:B:179:ARG:HG3	2.11	0.51
1:B:116:THR:HB	1:B:117:PRO:HD3	1.93	0.51
1:A:90:LEU:HD13	1:A:109:PRO:HA	1.93	0.50
1:B:329:ASN:HD22	1:B:364:ASN:HB2	1.76	0.50
3:B:503:CHD:H12	3:B:503:CHD:H212	1.93	0.50
1:B:212:GLN:HG3	1:B:213:VAL:N	2.27	0.50
1:A:102:PRO:O	1:A:103:ILE:C	2.55	0.50
1:A:412:ARG:HG3	1:A:415:LYS:HZ1	1.77	0.49
1:A:232:LEU:HD22	1:A:376:LEU:HD11	1.94	0.49
1:A:100:THR:O	1:A:101:LEU:HD23	2.13	0.49
1:A:273:GLY:O	1:B:298:ARG:HD2	2.12	0.49
3:A:501:CHD:H212	3:A:502:CHD:H193	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:GLU:HG3	1:A:159:TYR:CE2	2.48	0.48
1:A:226:ARG:HD3	1:A:279:GLU:OE1	2.13	0.48
1:B:171:GLU:CD	1:B:171:GLU:H	2.20	0.48
1:A:341:HIS:CG	1:A:342:ILE:N	2.79	0.48
3:A:503:CHD:H212	3:A:503:CHD:H12	1.95	0.48
1:A:175:GLU:O	1:A:179:ARG:HG3	2.14	0.47
1:B:374:ASN:ND2	1:B:374:ASN:C	2.72	0.47
1:B:400:THR:CA	1:B:415:LYS:HD2	2.41	0.47
1:B:304:LYS:HG3	1:B:310:TRP:CG	2.49	0.47
1:B:189:THR:HG22	1:B:224:ILE:O	2.14	0.47
1:B:337:PHE:CG	1:B:341:HIS:HE1	2.26	0.47
1:A:171:GLU:H	1:A:171:GLU:CD	2.21	0.47
1:A:213:VAL:HG13	1:A:215:ARG:CG	2.40	0.46
1:A:316:ASP:OD2	1:A:317:GLU:N	2.47	0.46
1:A:388:HIS:HD2	1:A:393:GLU:CD	2.23	0.46
1:A:71:ILE:N	1:A:71:ILE:HD12	2.31	0.46
1:A:210:TYR:HA	1:A:213:VAL:HG12	1.98	0.46
3:A:502:CHD:H212	3:A:502:CHD:H12	1.97	0.45
1:A:212:GLN:HG3	1:A:213:VAL:N	2.30	0.45
1:A:415:LYS:HB3	1:A:415:LYS:HZ2	1.81	0.45
1:A:193:GLN:OE1	1:A:336:ALA:HA	2.17	0.45
1:A:103:ILE:HG13	1:A:107:LEU:HG	1.99	0.45
1:A:158:LYS:HD3	1:A:160:TYR:OH	2.17	0.45
1:B:193:GLN:HG2	1:B:280:VAL:HA	1.98	0.45
1:A:253:ARG:NH1	4:A:575:HOH:O	2.51	0.44
1:A:298:ARG:HD2	1:B:273:GLY:O	2.17	0.44
1:B:374:ASN:HD22	1:B:375:PRO:CD	2.30	0.44
1:B:386:HIS:O	1:B:390:GLN:HG2	2.17	0.44
1:B:388:HIS:HD2	1:B:393:GLU:CD	2.26	0.44
1:A:357:ALA:HB1	1:A:362:VAL:CG1	2.46	0.44
1:A:96:ARG:NE	1:A:104:GLN:CD	2.76	0.44
1:A:67:PRO:HA	1:A:156:PRO:HG2	1.99	0.43
1:B:349:ASP:O	1:B:354:GLN:HG3	2.19	0.43
1:B:305:VAL:HG11	3:B:501:CHD:H112	2.00	0.43
1:A:250:LEU:HD12	1:A:250:LEU:N	2.21	0.43
1:B:216:LYS:HG3	1:B:217:PRO:HD2	2.00	0.43
1:A:251:GLU:H	1:A:251:GLU:CD	2.27	0.43
3:A:501:CHD:C21	3:A:502:CHD:H193	2.48	0.43
1:A:101:LEU:HB2	1:A:104:GLN:HG3	2.01	0.42
1:A:222:SER:OG	1:A:388:HIS:CE1	2.63	0.42
1:B:272:ARG:HH11	1:B:272:ARG:HG3	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:LEU:C	1:A:92:LEU:HD23	2.45	0.42
1:A:264:SER:HB2	1:A:301:TRP:HB3	2.02	0.42
1:A:263:CYS:HG	1:A:337:PHE:HZ	1.65	0.42
1:B:210:TYR:HA	1:B:213:VAL:HG12	2.01	0.42
1:B:332:LEU:HD13	1:B:348:LEU:HD22	2.02	0.42
3:A:501:CHD:H211	3:A:501:CHD:H232	1.85	0.42
1:B:349:ASP:O	1:B:353:SER:HB2	2.20	0.42
1:A:348:LEU:HD23	1:A:352:TYR:HD2	1.85	0.41
1:B:195:SER:HB3	1:B:198:THR:OG1	2.20	0.41
1:B:135:TRP:O	1:B:139:GLN:HG3	2.20	0.41
1:A:122:GLN:HG2	1:A:346:TYR:CE2	2.55	0.41
1:A:145:LYS:HB3	1:A:145:LYS:NZ	2.36	0.41
1:A:412:ARG:HA	1:A:415:LYS:NZ	2.35	0.41
1:A:177:MET:HB3	1:A:182:LEU:HD12	2.03	0.41
1:A:195:SER:HB3	1:A:198:THR:OG1	2.21	0.41
1:A:92:LEU:HD11	3:A:501:CHD:H7	2.01	0.41
1:A:288:MET:HG3	1:A:297:TYR:CD2	2.54	0.41
1:B:208:ARG:NH2	1:B:410:VAL:HG21	2.35	0.41
1:A:135:TRP:O	1:A:139:GLN:HG3	2.21	0.41
1:B:323:CYS:HG	1:B:362:VAL:HG12	1.84	0.41
1:B:314:GLN:HG2	4:B:1102:HOH:O	2.21	0.40
1:A:388:HIS:HD2	1:A:393:GLU:OE2	2.03	0.40
1:B:92:LEU:HD11	3:B:501:CHD:H152	2.03	0.40
3:B:503:CHD:O7	3:B:503:CHD:H41	2.21	0.40
1:A:313:PRO:HB3	1:B:271:ASN:O	2.21	0.40
1:A:108:ALA:HB3	1:A:109:PRO:CD	2.52	0.40

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	357/359 (99%)	345 (97%)	11 (3%)	1 (0%)	36	42
1	B	357/359 (99%)	347 (97%)	8 (2%)	2 (1%)	21	23
All	All	714/718 (99%)	692 (97%)	19 (3%)	3 (0%)	30	34

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	372	ASN
1	A	372	ASN
1	B	355	VAL

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	316/324 (98%)	312 (99%)	4 (1%)	61	76
1	B	316/324 (98%)	309 (98%)	7 (2%)	45	61
All	All	632/648 (98%)	621 (98%)	11 (2%)	53	69

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	145	LYS
1	A	193	GLN
1	A	298	ARG
1	A	415	LYS
1	B	184	ARG
1	B	219	MET
1	B	250	LEU
1	B	251	GLU
1	B	298	ARG
1	B	341	HIS
1	B	415	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	120	GLN
1	A	212	GLN
1	A	240	HIS
1	A	329	ASN
1	A	341	HIS
1	A	354	GLN
1	A	364	ASN
1	A	374	ASN
1	A	388	HIS
1	A	398	GLN
1	A	421	GLN
1	B	120	GLN
1	B	190	GLN
1	B	212	GLN
1	B	235	GLN
1	B	329	ASN
1	B	341	HIS
1	B	354	GLN
1	B	364	ASN
1	B	374	ASN
1	B	388	HIS
1	B	398	GLN
1	B	421	GLN

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](#)

8 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	FES	A	499	1	0,4,4	-	-	-		
3	CHD	B	501	-	32,32,32	0.81	2 (6%)	51,51,51	1.27	4 (7%)
3	CHD	B	502	-	32,32,32	0.85	2 (6%)	51,51,51	1.20	6 (11%)
3	CHD	A	501	-	32,32,32	0.83	2 (6%)	51,51,51	1.29	5 (9%)
3	CHD	A	503	-	32,32,32	0.81	2 (6%)	51,51,51	1.10	4 (7%)
2	FES	B	999	1	0,4,4	-	-	-		
3	CHD	B	503	-	32,32,32	0.81	1 (3%)	51,51,51	1.13	5 (9%)
3	CHD	A	502	-	32,32,32	0.83	1 (3%)	51,51,51	1.15	4 (7%)

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	FES	A	499	1	-	-	0/1/1/1
3	CHD	B	501	-	-	0/9/74/74	0/4/4/4
3	CHD	B	502	-	-	0/9/74/74	0/4/4/4
3	CHD	A	501	-	-	2/9/74/74	0/4/4/4
3	CHD	A	503	-	-	2/9/74/74	0/4/4/4
2	FES	B	999	1	-	-	0/1/1/1
3	CHD	B	503	-	-	0/9/74/74	0/4/4/4
3	CHD	A	502	-	-	0/9/74/74	0/4/4/4

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	CHD	C11-C12	2.77	1.57	1.53
3	A	502	CHD	C11-C12	2.58	1.57	1.53
3	B	501	CHD	C11-C12	2.31	1.57	1.53
3	A	501	CHD	C11-C12	2.30	1.57	1.53
3	B	503	CHD	C11-C12	2.29	1.56	1.53
3	A	503	CHD	C11-C12	2.22	1.56	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	CHD	O12-C12	-2.15	1.40	1.43
3	B	502	CHD	O12-C12	-2.11	1.40	1.43
3	A	503	CHD	O12-C12	-2.08	1.40	1.43
3	B	501	CHD	O12-C12	-2.04	1.40	1.43

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	CHD	C13-C17-C20	-3.78	114.91	119.48
3	B	501	CHD	C13-C17-C20	-3.62	115.10	119.48
3	A	503	CHD	C4-C5-C10	-3.20	109.26	112.66
3	B	503	CHD	C4-C5-C10	-2.98	109.48	112.66
3	B	502	CHD	C17-C13-C14	-2.89	97.21	100.11
3	A	501	CHD	C4-C5-C10	-2.83	109.65	112.66
3	B	501	CHD	C4-C5-C10	-2.80	109.68	112.66
3	B	502	CHD	C4-C5-C10	-2.73	109.76	112.66
3	B	501	CHD	C11-C12-C13	-2.66	108.56	111.26
3	A	501	CHD	C11-C12-C13	-2.59	108.63	111.26
3	A	502	CHD	C4-C5-C10	-2.53	109.96	112.66
3	B	503	CHD	C11-C12-C13	-2.41	108.81	111.26
3	B	502	CHD	C17-C13-C12	2.40	119.83	117.67
3	B	502	CHD	C1-C2-C3	-2.26	107.48	110.48
3	A	502	CHD	C11-C12-C13	-2.25	108.98	111.26
3	A	501	CHD	C1-C2-C3	-2.23	107.52	110.48
3	A	502	CHD	C1-C2-C3	-2.20	107.57	110.48
3	A	501	CHD	C21-C20-C22	-2.20	106.94	110.34
3	B	503	CHD	C13-C17-C20	-2.15	116.88	119.48
3	B	501	CHD	C1-C2-C3	-2.11	107.68	110.48
3	A	502	CHD	C6-C5-C10	-2.09	110.44	112.66
3	B	502	CHD	C11-C12-C13	-2.08	109.15	111.26
3	B	503	CHD	C1-C2-C3	-2.07	107.74	110.48
3	A	503	CHD	C17-C13-C14	-2.04	98.06	100.11
3	B	502	CHD	C6-C5-C10	-2.04	110.49	112.66
3	B	503	CHD	C11-C9-C10	-2.01	111.66	113.70
3	A	503	CHD	C11-C12-C13	-2.01	109.22	111.26
3	A	503	CHD	C6-C5-C10	-2.00	110.53	112.66

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	CHD	C22-C23-C24-O26

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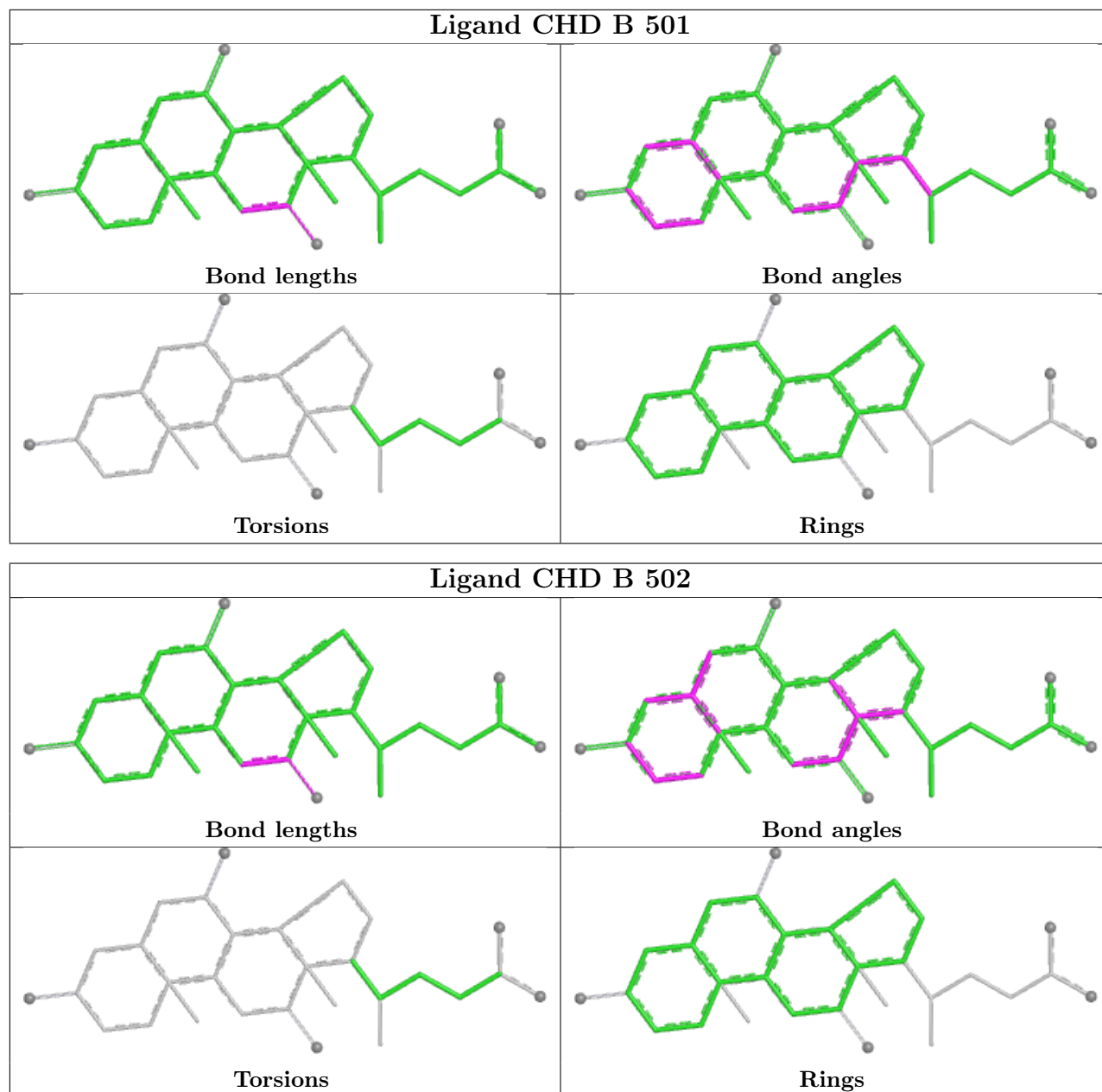
Mol	Chain	Res	Type	Atoms
3	A	501	CHD	C22-C23-C24-O25
3	A	503	CHD	C22-C23-C24-O25
3	A	503	CHD	C22-C23-C24-O26

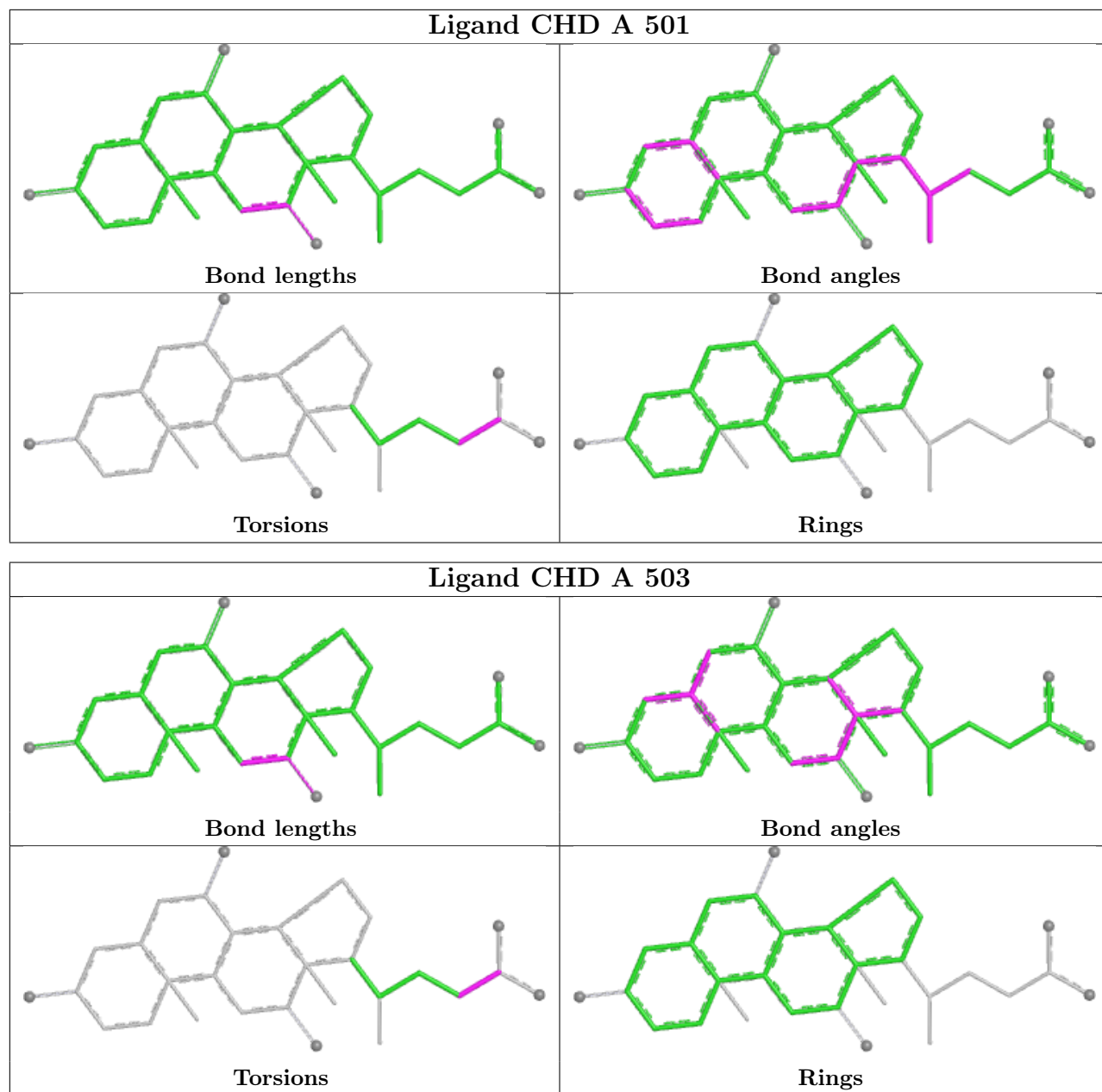
There are no ring outliers.

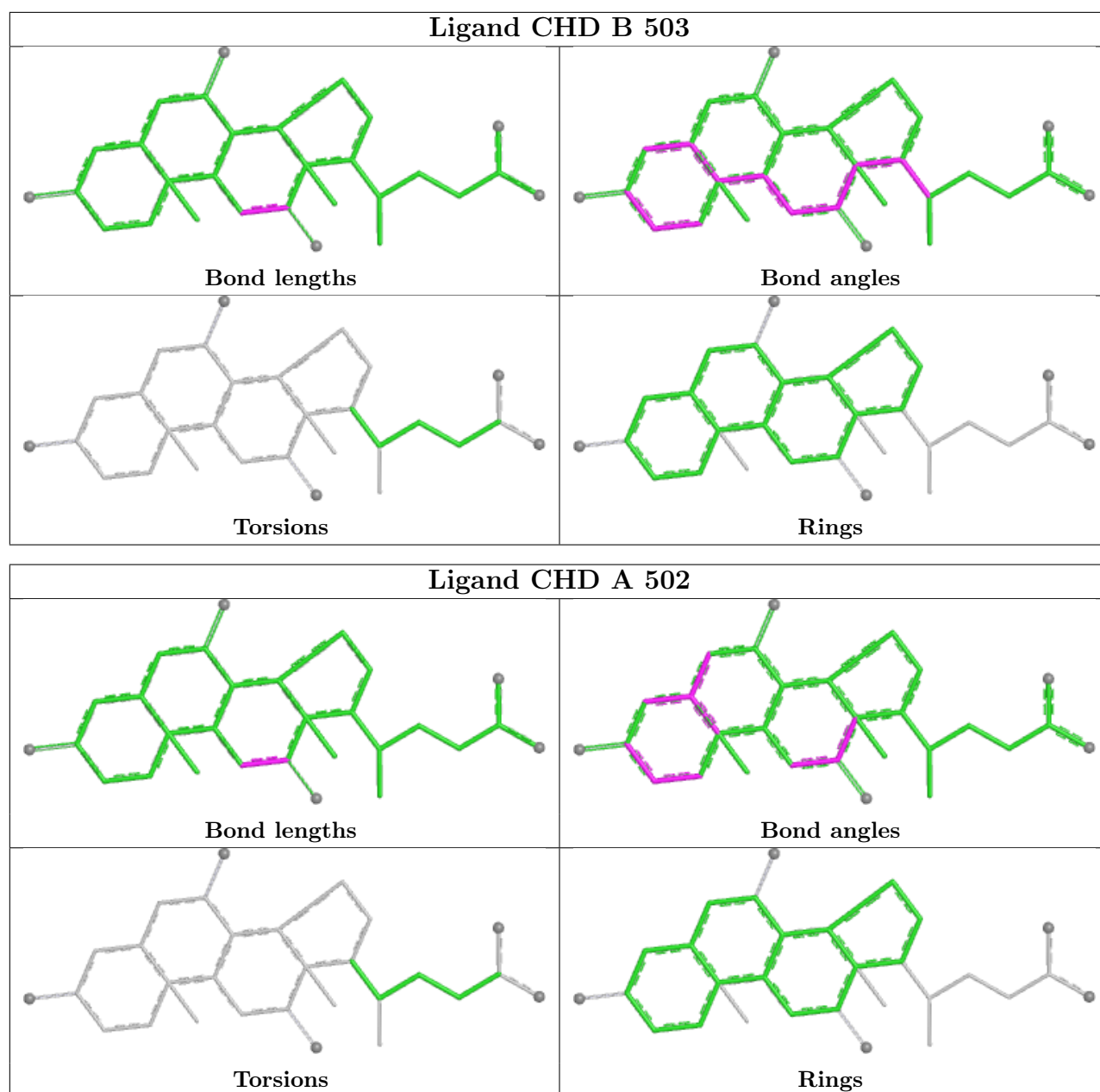
6 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	501	CHD	7	0
3	B	502	CHD	4	0
3	A	501	CHD	9	0
3	A	503	CHD	1	0
3	B	503	CHD	2	0
3	A	502	CHD	5	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

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.