



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 12:29 AM UTC

PDB ID : 1XUC / pdb_00001xuc
Title : Matrix metalloproteinase-13 complexed with non-zinc binding inhibitor
Authors : Engel, C.K.; Wendt, K.U.
Deposited on : 2004-10-26
Resolution : 1.70 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

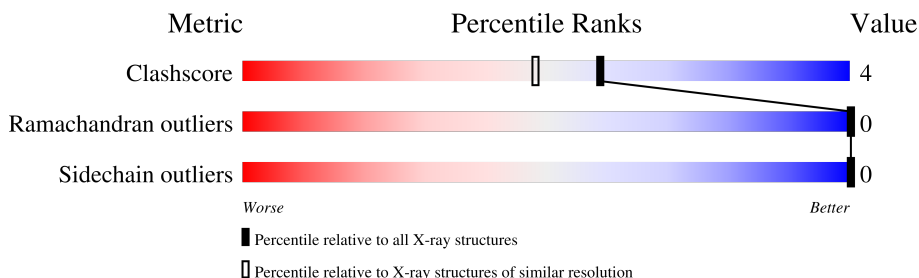
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)

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	171	 88% 11% ..
1	B	171	 84% 12% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3176 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 Collagenase 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	169	Total	C	N	O	S	0	0	0
			1346	870	217	254	5			
1	B	166	Total	C	N	O	S	0	0	0
			1321	857	214	245	5			

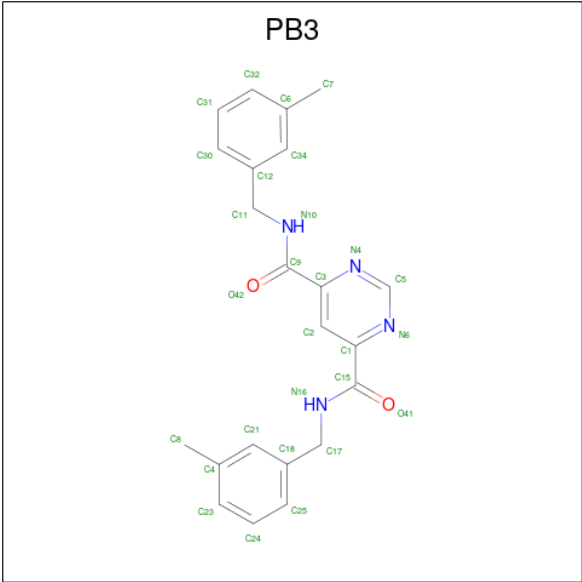
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Zn	0	0
			2	2		
2	B	2	Total	Zn	0	0
			2	2		

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 4 is N,N'-BIS(3-METHYLBENZYL)PYRIMIDINE-4,6-DICARBOXAMIDE (CCD ID: PB3) (formula: C₂₂H₂₂N₄O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			28	22	4	2		
4	B	1	Total	C	N	O	0	0
			28	22	4	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	243	Total	O	0	0
			243	243		
5	B	202	Total	O	0	0
			202	202		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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: Collagenase 3

Chain A:  88% 11% ..



- Molecule 1: Collagenase 3

Chain B:  84% 12% ..



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	134.44Å 36.48Å 95.32Å 90.00° 130.51° 90.00°	Depositor
Resolution (Å)	50.00 – 1.70	Depositor
% Data completeness (in resolution range)	98.2 (50.00-1.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.01	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.175 , 0.200	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3176	wwPDB-VP
Average B, all atoms (Å ²)	16.0	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: CA, ZN, PB3

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	0.32	0/1394	0.84	5/1895 (0.3%)
1	B	0.30	0/1369	0.88	11/1861 (0.6%)
All	All	0.31	0/2763	0.86	16/3756 (0.4%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	128	ASP	N-CA-C	7.38	120.38	111.82
1	B	128	ASP	N-CA-C	6.46	120.42	112.54
1	B	204	ASP	N-CA-C	-5.89	105.75	113.17
1	A	204	ASP	N-CA-C	-5.82	105.84	113.17
1	A	179	ASP	N-CA-C	5.59	119.28	111.56
1	B	179	ASP	N-CA-C	5.55	119.70	111.87
1	B	235	ASP	CA-C-N	5.37	125.69	119.47
1	B	235	ASP	C-N-CA	5.37	125.69	119.47
1	B	176	TYR	N-CA-C	-5.35	99.24	108.97
1	B	107	PHE	CA-C-N	5.31	125.23	119.76
1	B	107	PHE	C-N-CA	5.31	125.23	119.76
1	A	241	PHE	N-CA-C	-5.24	102.55	109.84
1	B	194	ASN	CB-CA-C	-5.08	110.32	117.23
1	B	241	PHE	N-CA-C	-5.07	103.09	110.08
1	A	243	ILE	N-CA-C	5.06	115.14	107.80
1	B	243	ILE	N-CA-C	5.01	115.19	108.17

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	1346	0	1250	14	0
1	B	1321	0	1236	13	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	28	0	22	0	0
4	B	28	0	22	0	0
5	A	243	0	0	2	0
5	B	202	0	0	3	0
All	All	3176	0	2530	23	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 4.

All (23) 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:109:ARG:HA	1:B:111:LEU:HD21	1.59	0.83
1:B:160:ILE:HG22	5:B:3138:HOH:O	1.93	0.68
1:A:107:PHE:HB3	1:B:109:ARG:HH21	1.59	0.67
1:A:133:GLU:HG2	5:A:2226:HOH:O	2.01	0.60
1:B:241:PHE:CD1	1:B:242:PRO:HD2	2.37	0.60
1:B:116:MET:HE3	1:B:150:PRO:HB2	1.88	0.55
1:A:111:LEU:HD23	1:A:190:PRO:HB2	1.90	0.54
1:A:241:PHE:CD1	1:A:242:PRO:HD2	2.43	0.54
1:B:129:MET:HE3	1:B:209:SER:OG	2.10	0.52
1:A:246:TYR:CE1	1:A:248:GLY:HA2	2.46	0.50
1:A:117:ASN:ND2	1:A:152:ASN:HD22	2.12	0.47
1:B:131:HIS:HD2	5:B:3030:HOH:O	1.96	0.47
1:B:246:TYR:CE1	1:B:248:GLY:HA2	2.52	0.45
1:B:182:SER:HA	5:B:3161:HOH:O	2.16	0.45
1:A:112:LYS:HE3	5:A:2221:HOH:O	2.17	0.45
1:B:181:PRO:O	1:B:182:SER:HB3	2.18	0.44
1:A:149:THR:HB	1:A:150:PRO:CD	2.48	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:THR:C	1:B:109:ARG:HD3	2.43	0.42
1:A:246:TYR:HE1	1:A:248:GLY:HA2	1.83	0.42
1:B:160:ILE:HG23	1:B:160:ILE:O	2.20	0.41
1:A:111:LEU:CD2	1:A:190:PRO:HB2	2.49	0.41
1:A:109:ARG:O	1:B:109:ARG:HB2	2.20	0.40
1:A:109:ARG:HH11	1:A:109:ARG:HG2	1.86	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	167/171 (98%)	160 (96%)	7 (4%)	0	100	100
1	B	164/171 (96%)	156 (95%)	8 (5%)	0	100	100
All	All	331/342 (97%)	316 (96%)	15 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	144/146 (99%)	144 (100%)	0	100	100
1	B	141/146 (97%)	141 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	285/292 (98%)	285 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	105	ASN
1	A	117	ASN
1	A	157	HIS
1	B	131	HIS
1	B	194	ASN
1	B	263	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 ⓘ

Of 10 ligands modelled in this entry, 8 are monoatomic - leaving 2 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
4	PB3	B	3001	-	30,30,30	2.39	18 (60%)	40,40,40	2.78	11 (27%)
4	PB3	A	2001	-	30,30,30	2.41	17 (56%)	40,40,40	2.80	11 (27%)

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
4	PB3	B	3001	-	-	0/18/18/18	0/3/3/3
4	PB3	A	2001	-	-	0/18/18/18	0/3/3/3

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	2001	PB3	C3-N4	4.17	1.42	1.34
4	B	3001	PB3	C3-N4	4.07	1.41	1.34
4	A	2001	PB3	C1-N6	3.94	1.41	1.34
4	B	3001	PB3	C1-N6	3.89	1.41	1.34
4	B	3001	PB3	C30-C12	3.70	1.46	1.38
4	A	2001	PB3	C25-C18	3.68	1.46	1.38
4	A	2001	PB3	C21-C4	3.67	1.44	1.39
4	A	2001	PB3	C5-N4	3.63	1.40	1.33
4	A	2001	PB3	C30-C12	3.56	1.45	1.38
4	B	3001	PB3	C25-C18	3.52	1.45	1.38
4	B	3001	PB3	C21-C4	3.51	1.44	1.39
4	B	3001	PB3	C5-N4	3.48	1.40	1.33
4	A	2001	PB3	C5-N6	3.33	1.39	1.33
4	B	3001	PB3	C5-N6	3.29	1.39	1.33
4	A	2001	PB3	C24-C25	2.98	1.44	1.38
4	B	3001	PB3	C34-C6	2.97	1.43	1.39
4	B	3001	PB3	C24-C25	2.88	1.43	1.38
4	B	3001	PB3	C15-N16	2.86	1.40	1.33
4	A	2001	PB3	C34-C6	2.78	1.43	1.39
4	A	2001	PB3	C15-N16	2.78	1.39	1.33
4	B	3001	PB3	C31-C30	2.40	1.43	1.38
4	A	2001	PB3	C24-C23	2.39	1.43	1.38
4	B	3001	PB3	C34-C12	2.39	1.43	1.39
4	A	2001	PB3	C21-C18	2.35	1.43	1.39
4	B	3001	PB3	C31-C32	2.32	1.42	1.38
4	B	3001	PB3	C21-C18	2.31	1.43	1.39
4	A	2001	PB3	C31-C30	2.30	1.42	1.38

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	2001	PB3	C34-C12	2.27	1.43	1.39
4	A	2001	PB3	C31-C32	2.26	1.42	1.38
4	A	2001	PB3	C23-C4	2.18	1.44	1.38
4	B	3001	PB3	C24-C23	2.17	1.42	1.38
4	B	3001	PB3	C9-N10	2.16	1.38	1.33
4	B	3001	PB3	C23-C4	2.13	1.43	1.38
4	B	3001	PB3	C32-C6	2.08	1.43	1.38
4	A	2001	PB3	C9-N10	2.03	1.38	1.33

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2001	PB3	C5-N6-C1	10.25	121.59	115.81
4	B	3001	PB3	C5-N6-C1	10.00	121.44	115.81
4	B	3001	PB3	C5-N4-C3	9.11	120.94	115.81
4	A	2001	PB3	C5-N4-C3	8.93	120.84	115.81
4	A	2001	PB3	N4-C5-N6	-7.46	117.29	128.58
4	B	3001	PB3	N4-C5-N6	-7.39	117.39	128.58
4	A	2001	PB3	C15-C1-N6	2.81	120.70	117.42
4	B	3001	PB3	O42-C9-C3	2.80	127.28	121.08
4	A	2001	PB3	O42-C9-C3	2.76	127.17	121.08
4	B	3001	PB3	C32-C6-C34	2.74	122.01	117.96
4	B	3001	PB3	C23-C4-C21	2.67	121.90	117.96
4	A	2001	PB3	C32-C6-C34	2.67	121.90	117.96
4	B	3001	PB3	C15-C1-N6	2.64	120.50	117.42
4	A	2001	PB3	C23-C4-C21	2.63	121.83	117.96
4	A	2001	PB3	O41-C15-C1	2.32	126.21	121.08
4	A	2001	PB3	O41-C15-N16	-2.32	118.07	122.59
4	B	3001	PB3	O41-C15-C1	2.31	126.18	121.08
4	A	2001	PB3	C9-C3-N4	2.25	120.05	117.42
4	B	3001	PB3	O41-C15-N16	-2.20	118.30	122.59
4	B	3001	PB3	C9-C3-N4	2.18	119.97	117.42
4	A	2001	PB3	C2-C1-N6	-2.10	120.75	123.27
4	B	3001	PB3	O42-C9-N10	-2.03	118.63	122.59

There are no chirality outliers.

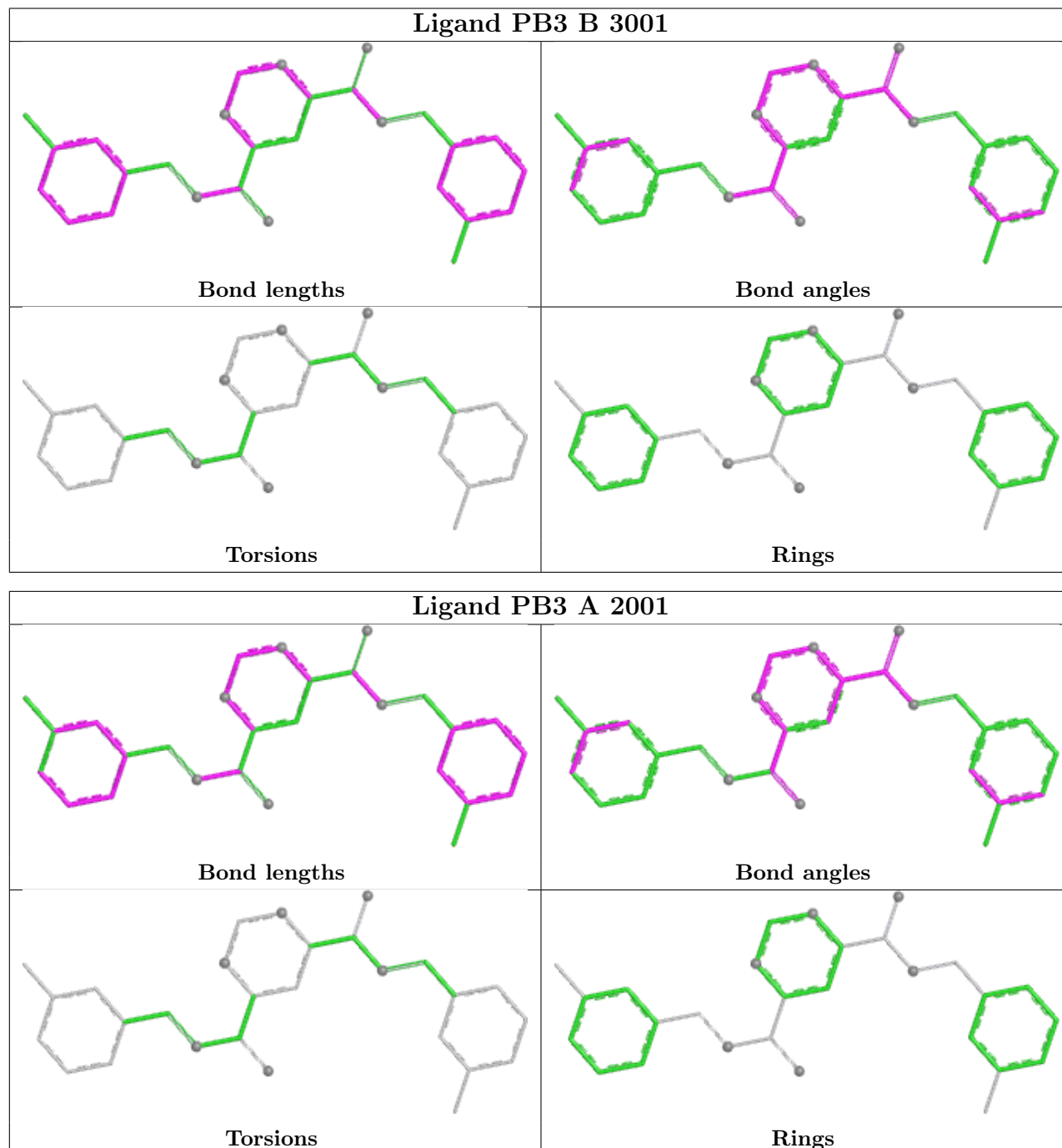
There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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.