



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 28, 2026 – 11:56 PM UTC

PDB ID : 1L7V / pdb_0000117v
Title : Bacterial ABC Transporter Involved in B12 Uptake
Authors : Locher, K.P.; Lee, A.T.; Rees, D.C.
Deposited on : 2002-03-18
Resolution : 3.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

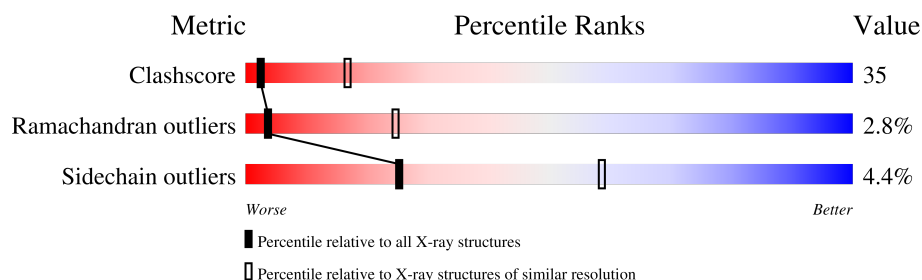
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)

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	326	
1	B	326	
2	C	249	
2	D	249	

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
3	V4O	C	250	-	-	X	-
3	V4O	D	250	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8410 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 VITAMIN B12 TRANSPORT SYSTEM PERMEASE PROTEIN BTUC.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	Se	163	0	0
			2441	1611	418	393	8	11			
1	B	324	Total	C	N	O	S	Se	183	0	0
			2441	1611	418	393	8	11			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP P06609
A	23	MSE	MET	modified residue	UNP P06609
A	77	MSE	MET	modified residue	UNP P06609
A	160	MSE	MET	modified residue	UNP P06609
A	176	MSE	MET	modified residue	UNP P06609
A	179	MSE	MET	modified residue	UNP P06609
A	180	MSE	MET	modified residue	UNP P06609
A	194	MSE	MET	modified residue	UNP P06609
A	211	MSE	MET	modified residue	UNP P06609
A	213	MSE	MET	modified residue	UNP P06609
A	242	MSE	MET	modified residue	UNP P06609
B	1	MSE	MET	modified residue	UNP P06609
B	23	MSE	MET	modified residue	UNP P06609
B	77	MSE	MET	modified residue	UNP P06609
B	160	MSE	MET	modified residue	UNP P06609
B	176	MSE	MET	modified residue	UNP P06609
B	179	MSE	MET	modified residue	UNP P06609
B	180	MSE	MET	modified residue	UNP P06609
B	194	MSE	MET	modified residue	UNP P06609
B	211	MSE	MET	modified residue	UNP P06609
B	213	MSE	MET	modified residue	UNP P06609
B	242	MSE	MET	modified residue	UNP P06609

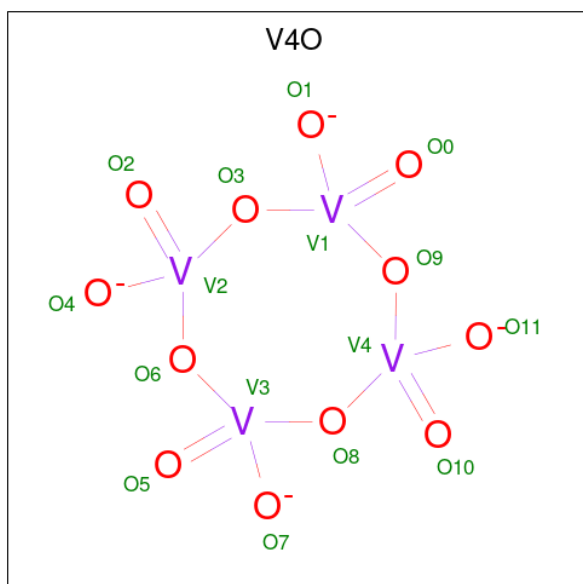
- Molecule 2 is a protein called Vitamin B12 transport ATP-binding protein btuD.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	C	231	Total	C	N	O	S	Se	59	0	0
			1748	1094	322	324	1	7			
2	D	231	Total	C	N	O	S	Se	59	0	0
			1748	1094	322	324	1	7			

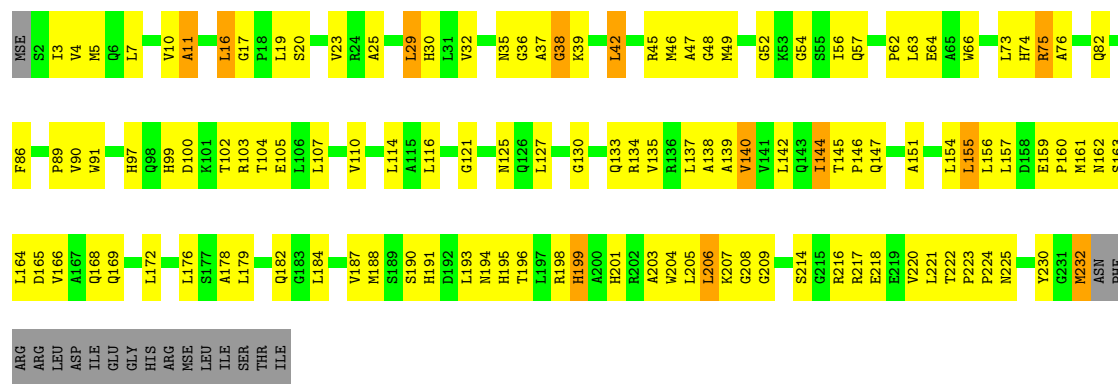
There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	MSE	MET	modified residue	UNP P06611
C	5	MSE	MET	modified residue	UNP P06611
C	46	MSE	MET	modified residue	UNP P06611
C	49	MSE	MET	modified residue	UNP P06611
C	161	MSE	MET	modified residue	UNP P06611
C	188	MSE	MET	modified residue	UNP P06611
C	211	MSE	MET	modified residue	UNP P06611
C	232	MSE	MET	modified residue	UNP P06611
C	244	MSE	MET	modified residue	UNP P06611
D	1	MSE	MET	modified residue	UNP P06611
D	5	MSE	MET	modified residue	UNP P06611
D	46	MSE	MET	modified residue	UNP P06611
D	49	MSE	MET	modified residue	UNP P06611
D	161	MSE	MET	modified residue	UNP P06611
D	188	MSE	MET	modified residue	UNP P06611
D	211	MSE	MET	modified residue	UNP P06611
D	232	MSE	MET	modified residue	UNP P06611
D	244	MSE	MET	modified residue	UNP P06611

- Molecule 3 is CYCLO-TETRAMETAVANADATE (CCD ID: V4O) (formula: $O_{12}V_4$).

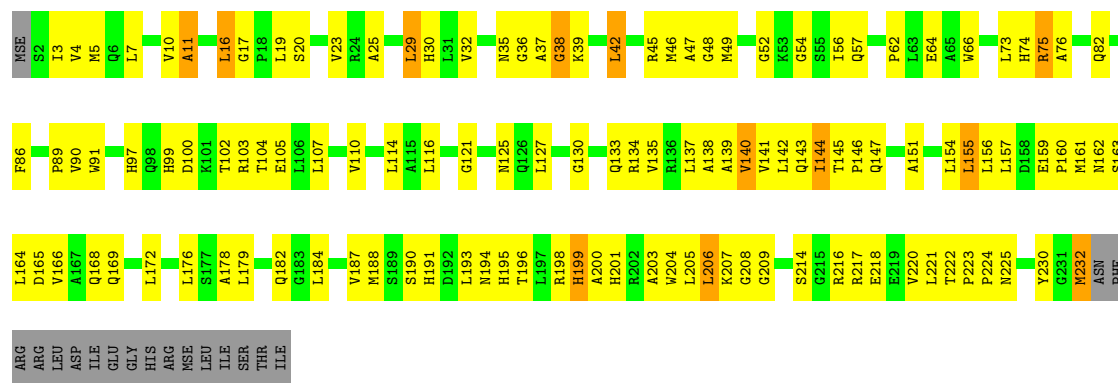


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	O	V	0	0
			16	12	4		
3	D	1	Total	O	V	0	0
			16	12	4		



● Molecule 2: Vitamin B12 transport ATP-binding protein btuD

Chain D: 43% 45% 5% 7%



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	92.99Å 106.29Å 95.54Å 90.00° 99.29° 90.00°	Depositor
Resolution (Å)	15.00 – 3.20	Depositor
% Data completeness (in resolution range)	(Not available) (15.00-3.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.262 , 0.286	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8410	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: V4O

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	0.43	0/2485	0.92	8/3378 (0.2%)
1	B	0.46	0/2485	0.95	11/3378 (0.3%)
2	C	0.41	0/1773	0.90	4/2395 (0.2%)
2	D	0.39	0/1773	0.89	4/2395 (0.2%)
All	All	0.43	0/8516	0.92	27/11546 (0.2%)

There are no bond length outliers.

The worst 5 of 27 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	36	GLN	OE1-CD-NE2	-9.84	112.76	122.60
1	A	36	GLN	OE1-CD-NE2	-9.12	113.48	122.60
1	A	256	ILE	N-CA-C	-8.77	104.94	113.53
1	B	259	VAL	N-CA-C	8.39	118.31	110.42
1	B	83	ASN	CA-C-N	8.13	128.22	119.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2441	0	2606	178	1
1	B	2441	0	2606	164	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	1748	0	1775	120	0
2	D	1748	0	1775	122	0
3	C	16	0	0	6	0
3	D	16	0	0	6	0
All	All	8410	0	8762	561	1

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 35.

The worst 5 of 561 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)
2:C:30:HIS:HE2	2:C:188:MSE:HE3	1.02	1.17
2:D:30:HIS:HE2	2:D:188:MSE:HE3	1.02	1.09
1:A:77:MSE:HE1	1:A:88:PRO:HG3	1.36	1.07
1:B:77:MSE:HE1	1:B:88:PRO:HG3	1.35	1.05
1:A:269:LEU:HD22	1:A:274:VAL:HG11	1.41	1.02

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:ARG:NH1	1:B:35:GLU:OE2[2_657]	1.49	0.71

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	322/326 (99%)	258 (80%)	53 (16%)	11 (3%)	3	20
1	B	322/326 (99%)	255 (79%)	55 (17%)	12 (4%)	2	18

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	229/249 (92%)	194 (85%)	31 (14%)	4 (2%)	7	35
2	D	229/249 (92%)	195 (85%)	30 (13%)	4 (2%)	7	35
All	All	1102/1150 (96%)	902 (82%)	169 (15%)	31 (3%)	4	25

5 of 31 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	51	PHE
1	B	181	GLY
2	C	103	ARG
2	D	103	ARG
1	A	172	LEU

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	249/239 (104%)	235 (94%)	14 (6%)	19	52
1	B	249/239 (104%)	241 (97%)	8 (3%)	34	65
2	C	184/192 (96%)	176 (96%)	8 (4%)	26	59
2	D	184/192 (96%)	176 (96%)	8 (4%)	26	59
All	All	866/862 (100%)	828 (96%)	38 (4%)	25	59

5 of 38 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	C	206	LEU
2	D	155	LEU
2	C	232	MSE
2	D	42	LEU
2	D	232	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 28 such sidechains are listed below:

Mol	Chain	Res	Type
2	C	194	ASN
2	D	228	GLN
2	D	35	ASN
2	D	168	GLN
2	C	228	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](#)

2 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	V4O	C	250	-	12,16,16	1.69	3 (25%)	-		
3	V4O	D	250	-	12,16,16	1.72	3 (25%)	-		

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	V4O	C	250	-	-	-	0/1/1/1
3	V4O	D	250	-	-	-	0/1/1/1

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	250	V4O	O0-V1	-3.45	1.58	1.62
3	C	250	V4O	O0-V1	-3.37	1.58	1.62
3	C	250	V4O	O8-V4	-2.46	1.71	1.77
3	D	250	V4O	O8-V4	-2.42	1.71	1.77
3	C	250	V4O	O3-V2	-2.37	1.71	1.77

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

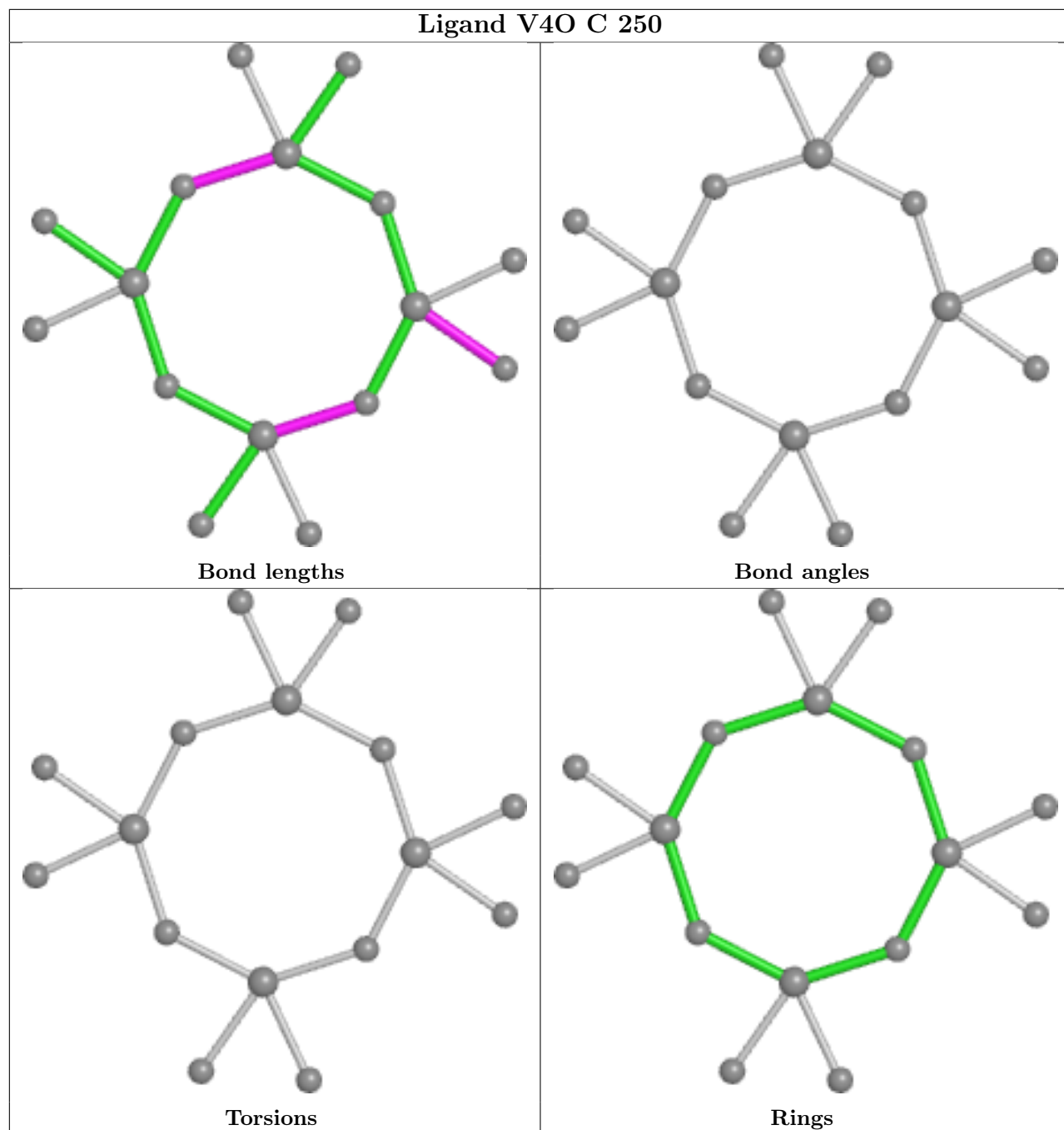
There are no ring outliers.

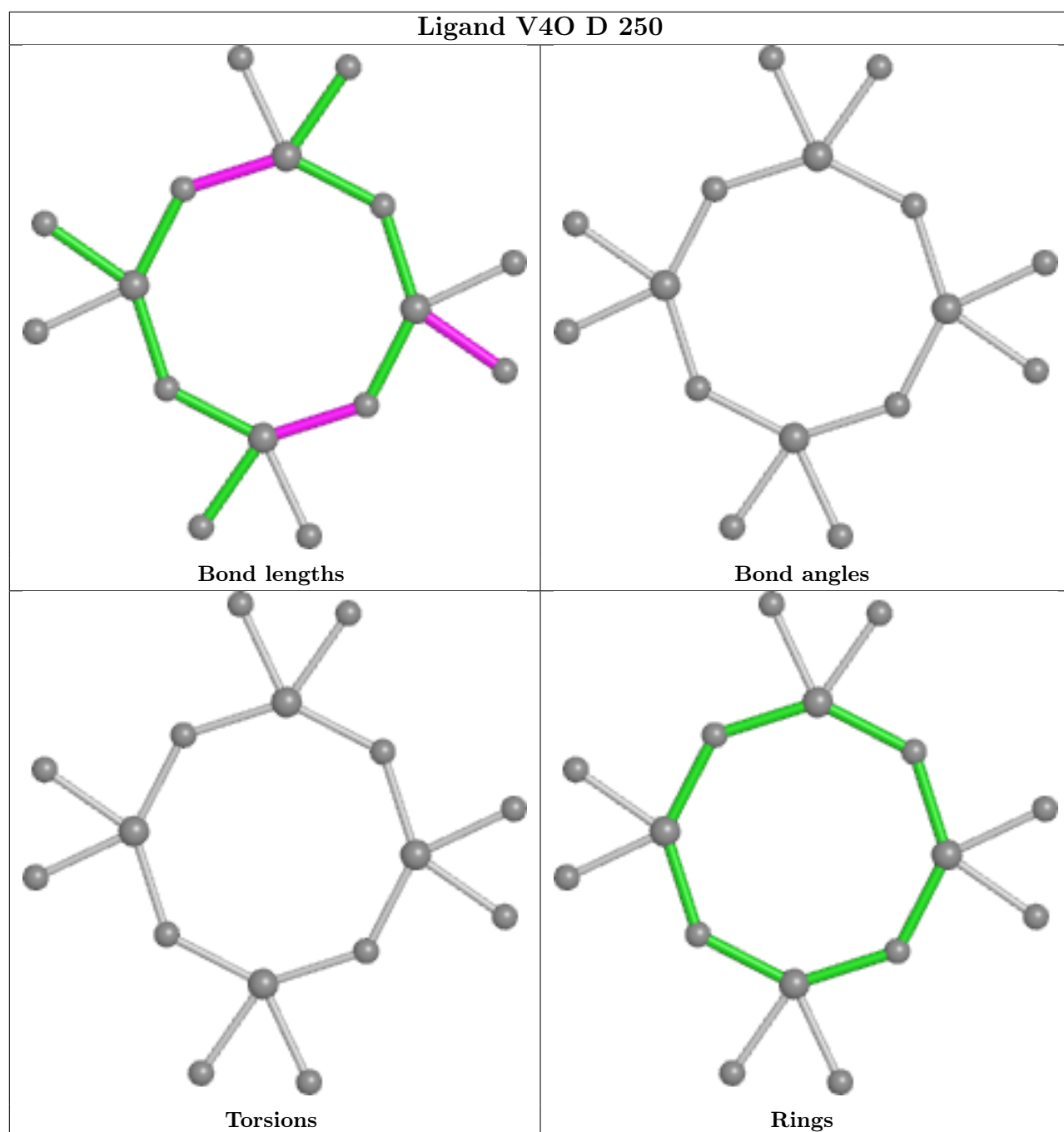
2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	250	V4O	6	0
3	D	250	V4O	6	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand V4O C 250





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.