

wwPDB X-ray Structure Validation Summary Report (i)

Sep 14, 2023 – 06:30 AM EDT

PDB ID : 1QGW

Title : CRYSTAL STRUCTURE OF PHYCOERYTHRIN 545 FROM THE MA-

RINE CRYPTOPHYTE RHODOMONAS CS24

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Deposited on : 1999-05-10

Resolution : 1.63 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 (i) 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 (1)) were used in the production of this report:

MolProbity: 4.02b-467

Mogul : 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : NOT EXECUTED

EDS : NOT EXECUTED

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

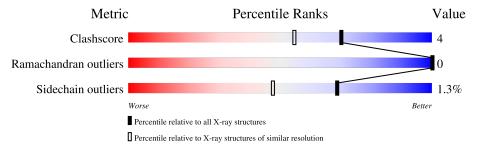
Validation Pipeline (wwPDB-VP) : 2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\ DIFFRACTION$

The reported resolution of this entry is 1.63 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	Similar resolution		
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$		
Clashscore	141614	3268 (1.66-1.62)		
Ramachandran outliers	138981	3215 (1.66-1.62)		
Sidechain outliers	138945	3215 (1.66-1.62)		

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain	
1	A	76	95%	5%
2	В	67	93%	7%
3	С	177	90%	5% • 5%
3	D	177	94%	5% •



2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 4788 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 PROTEIN (CRYPTOPHYTAN PHYCOERYTHRIN (ALPHA-1 CHAIN)).

Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf	Trace
1	A	76	Total 567	C 348	N 98	O 117	S 4	6	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Α	4	LYZ	LYS	modified residue	UNP Q00433
A	10	GLN	LEU	conflict	UNP Q00433

• Molecule 2 is a protein called PROTEIN (CRYPTOPHYTAN PHYCOERYTHRIN (ALPHA-2 CHAIN)).

Mol	Chain	Residues		Atoms				ZeroOcc	AltConf	Trace
2	В	67	Total 493	C 303	N 86	O 98	S 6	4	2	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
В	4	LYZ	LYS	modified residue	UNP Q00433

• Molecule 3 is a protein called PROTEIN (CRYPTOPHYTAN PHYCOERYTHRIN (BETA CHAIN)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	С	169	Total	С	N	О	S	28	4	0
ა		109	1236	759	215	252	10	20	4	U
2	D	177	Total	С	N	О	S	19	0	0
3	ט	111	1300	794	225	270	11	19	9	U

There are 12 discrepancies between the modelled and reference sequences:

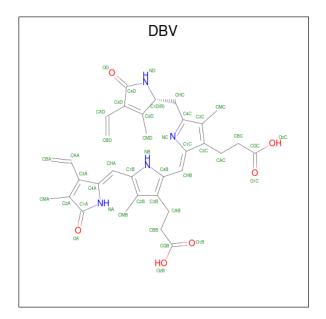


Chain	Residue	Modelled	Actual	Comment	Reference
С	50	CYS	VAL	conflict	UNP P27198
С	56	VAL	TYR	conflict	UNP P27198
С	61	CYS	GLU	conflict	UNP P27198
С	65	SER	HIS	conflict	UNP P27198
С	72	MEN	ASN	modified residue	UNP P27198
С	73	CYS	GLU	conflict	UNP P27198
D	50	CYS	VAL	conflict	UNP P27198
D	56	VAL	TYR	conflict	UNP P27198
D	61	CYS	GLU	conflict	UNP P27198
D	65	SER	HIS	conflict	UNP P27198
D	72	MEN	ASN	modified residue	UNP P27198
D	73	CYS	GLU	conflict	UNP P27198

• Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Mg 1 1	0	0
4	D	1	Total Mg 1 1	0	0

 $\bullet \ \ Molecule\ 5\ is\ 15,16-DIHYDROBILIVERDIN\ (three-letter\ code:\ DBV)\ (formula:\ C_{33}H_{36}N_4O_6).$



Mol	Chain	Residues	A	${f Atoms}$			ZeroOcc	AltConf
5	A	1	Total 43	C 33	N 4	O 6	0	0

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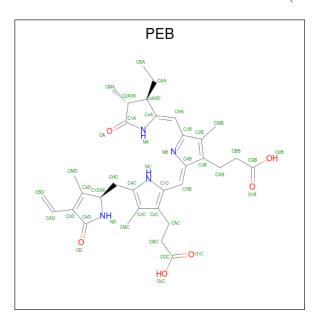
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Mol	Chain	Residues	A	Aton	ns		ZeroOcc	AltConf
5	В	1	Total 43	C 33	N 4	O 6	0	0

• Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	С	1	Total Cl 1 1	0	0

 \bullet Molecule 7 is PHYCOERYTHROBILIN (three-letter code: PEB) (formula: $C_{33}H_{40}N_4O_6).$



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	С	1	Total	С	N	О	0	0
'		1	43	33	4	6	U	
7	С	1	Total	С	N	О	0	0
•		1	43	33	4	6	U	0
7	C	1	Total	С	N	Ο	0	0
1		1	43	33	4	6	U	0
7	D	1	Total	\mathbf{C}	N	Ο	0	0
•	D	1	43	33	4	6	U	
7	D	1	Total	\mathbf{C}	N	Ο	0	0
•	D	1	43	33	4	6	O	
7	D	1	Total	С	N	O	0	0
'	D	1	43	33	4	6		

• Molecule 8 is water.



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	168	Total O 168 168	0	0
8	В	135	Total O 135 135	0	0
8	С	217	Total O 217 217	0	0
8	D	325	Total O 325 325	0	0

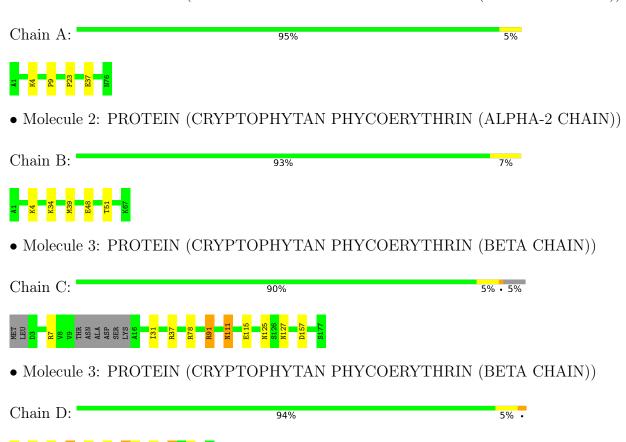


3 Residue-property plots (i)

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.







4 Data and refinement statistics (i)

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

Property	Value	Source	
Space group	P 21 21 21	Depositor	
Cell constants	63.03Å 82.63Å 89.55Å	Depositor	
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	20.00 - 1.63	Depositor	
% Data completeness	95.4 (20.00-1.63)	Depositor	
(in resolution range)	30.4 (20.00 1.00)		
R_{merge}	0.04	Depositor	
R_{sym}	0.04	Depositor	
Refinement program	REFMAC	Depositor	
R, R_{free}	0.149 , 0.188	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	4788	wwPDB-VP	
Average B, all atoms (Å ²)	8.0	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: MEN, PEB, MG, DBV, CL, LYZ

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).

Mal Chain		Bo	nd lengths	Bond angles		
Mol	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.44	0/563	0.86	0/749	
2	В	0.41	0/498	0.81	0/661	
3	С	0.44	0/1258	0.90	3/1695 (0.2%)	
3	D	0.45	1/1348 (0.1%)	0.88	5/1814 (0.3%)	
All	All	0.44	$1/3667 \ (0.0\%)$	0.88	8/4919 (0.2%)	

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	Observed(A)	$\operatorname{Ideal}(ext{\AA})$
3	D	7	ARG	CD-NE	-5.34	1.37	1.46

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
3	С	91	ARG	NE-CZ-NH2	-10.33	115.13	120.30
3	D	108	ARG	NE-CZ-NH2	7.33	123.96	120.30
3	С	7	ARG	NE-CZ-NH2	7.29	123.94	120.30
3	D	1	MET	CG-SD-CE	5.97	109.75	100.20
3	D	7	ARG	CD-NE-CZ	5.55	131.37	123.60

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	567	0	564	3	0
2	В	493	0	511	5	2
3	С	1236	0	1231	5	2
3	D	1300	0	1300	5	0
4	A	1	0	0	0	0
4	D	1	0	0	0	0
5	A	43	0	33	3	0
5	В	43	0	33	4	0
6	С	1	0	0	0	0
7	С	129	0	110	4	0
7	D	129	0	110	3	0
8	A	168	0	0	1	1
8	В	135	0	0	1	1
8	С	217	0	0	2	4
8	D	325	0	0	2	4
All	All	4788	0	3892	28	7

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.

The worst 5 of 28 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
3:C:115:GLU:HG3	8:C:1103:HOH:O	1.76	0.85
2:B:51:THR:HG21	8:D:1087:HOH:O	1.82	0.80
1:A:37:GLU:OE2	8:A:929:HOH:O	2.09	0.70
1:A:9:PRO:O	3:C:91:ARG:HD3	1.94	0.68
3:C:125:ASN:OD1	8:C:1115:HOH:O	2.14	0.64

The worst 5 of 7 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	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
2:B:34:LYS:NZ	3:C:157:ASP:OD1[4_456]	1.73	0.47
8:A:1028:HOH:O	8:D:1103:HOH:O[3_555]	1.91	0.29
8:C:1047:HOH:O	8:D:912:HOH:O[3_555]	2.03	0.17
2:B:34:LYS:NZ	3:C:157:ASP:CG[4_456]	2.10	0.10
8:B:302:HOH:O	8:C:1097:HOH:O[1_455]	2.10	0.10



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	Perce	ntiles
1	A	73/76 (96%)	72 (99%)	1 (1%)	0	100	100
2	В	66/67 (98%)	64~(97%)	2 (3%)	0	100	100
3	\mathbf{C}	168/177~(95%)	167 (99%)	1 (1%)	0	100	100
3	D	182/177 (103%)	181 (100%)	1 (0%)	0	100	100
All	All	489/497 (98%)	484 (99%)	5 (1%)	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	Perce	ntiles
1	A	60/60 (100%)	60 (100%)	0	100	100
2	В	52/50 (104%)	52 (100%)	0	100	100
3	С	138/141 (98%)	136 (99%)	2 (1%)	67	45
3	D	150/141 (106%)	147 (98%)	3 (2%)	55	29
All	All	400/392 (102%)	395 (99%)	5 (1%)	69	47

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	С	111	ASN
3	С	127	ASN

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Mol	Chain	Res	Type
3	D	111	ASN
3	D	127	ASN
3	D	144	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
3	С	111	ASN
3	С	127	ASN
3	D	111	ASN
3	D	127	ASN
3	D	144	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)

4 non-standard protein/DNA/RNA residues 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	Tuno	Chain	Res	Link	B	ond leng	gths	Bond angles		
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
2	LYZ	В	4	2	7,9,10	0.71	0	4,10,12	3.83	2 (50%)
3	MEN	С	72	3	7,8,9	0.56	0	6,9,11	0.49	0
3	MEN	D	72	3	7,8,9	0.40	0	6,9,11	0.48	0
1	LYZ	A	4	1	7,9,10	0.65	0	4,10,12	2.20	2 (50%)

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	LYZ	В	4	2	-	1/8/9/11	-
3	MEN	С	72	3	-	2/7/8/10	_
3	MEN	D	72	3	-	2/7/8/10	-
1	LYZ	A	4	1	-	1/8/9/11	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\mathbf{Ideal}(^{o})$
2	В	4	LYZ	OH-CD-CE	-6.24	88.08	109.33
2	В	4	LYZ	OH-CD-CG	-4.03	97.69	109.21
1	A	4	LYZ	OH-CD-CG	-3.70	98.63	109.21
1	A	4	LYZ	OH-CD-CE	-2.26	101.63	109.33

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	4	LYZ	O-C-CA-CB
2	В	4	LYZ	CG-CD-CE-NZ
3	D	72	MEN	CA-CB-CG-OD1
3	D	72	MEN	CA-CB-CG-ND2
3	С	72	MEN	CA-CB-CG-OD1

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 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	Trno	Chain	Res	Link	Вс	ond leng	ths	В	ond ang	gles
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	PEB	С	250	3	43,46,46	1.79	7 (16%)	45,67,67	2.01	13 (28%)
7	PEB	D	258	3	43,46,46	1.72	5 (11%)	45,67,67	2.11	10 (22%)
7	PEB	D	282	3	43,46,46	1.88	6 (13%)	45,67,67	2.00	10 (22%)
7	PEB	С	258	3	43,46,46	1.80	6 (13%)	45,67,67	2.11	7 (15%)
5	DBV	В	219	2	42,46,46	1.69	5 (11%)	42,67,67	1.60	8 (19%)
7	PEB	С	282	3	43,46,46	1.84	7 (16%)	45,67,67	2.02	12 (26%)
7	PEB	D	250	3	43,46,46	1.89	7 (16%)	45,67,67	2.02	15 (33%)
5	DBV	A	219	1	42,46,46	1.81	6 (14%)	42,67,67	1.54	8 (19%)

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
7	PEB	С	250	3	-	5/24/74/74	0/4/4/4
7	PEB	D	258	3	-	5/24/74/74	0/4/4/4
7	PEB	D	282	3	-	2/24/74/74	0/4/4/4
7	PEB	С	258	3	-	5/24/74/74	0/4/4/4
5	DBV	В	219	2	-	5/26/74/74	0/4/4/4
7	PEB	С	282	3	-	5/24/74/74	0/4/4/4
7	PEB	D	250	3	-	8/24/74/74	0/4/4/4
5	DBV	A	219	1	-	7/26/74/74	0/4/4/4

The worst 5 of 49 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\mathring{\mathrm{A}})$	Ideal(Å)
7	D	282	PEB	C2D-C3D	7.41	1.44	1.34
5	A	219	DBV	CHB-C1C	7.26	1.41	1.35
7	С	282	PEB	C2D-C3D	7.16	1.43	1.34
7	D	258	PEB	C2D-C3D	6.85	1.43	1.34
7	D	250	PEB	C2A-C1A	-6.68	1.46	1.52

The worst 5 of 83 bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$Ideal(^{o})$
7	D	258	PEB	C1C-CHB-C4B	10.06	140.83	128.81
7	С	258	PEB	C1C-CHB-C4B	9.92	140.66	128.81
7	D	282	PEB	C1C-CHB-C4B	8.65	139.14	128.81
7	С	282	PEB	C1C-CHB-C4B	8.23	138.64	128.81
7	С	250	PEB	C1C-CHB-C4B	7.69	137.99	128.81

There are no chirality outliers.

5 of 42 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	219	DBV	C2A-C3A-CAA-CBA
5	A	219	DBV	C4A-C3A-CAA-CBA
5	A	219	DBV	NB-C1B-CHA-C4A
5	A	219	DBV	C2B-C1B-CHA-C4A
5	В	219	DBV	C4A-C3A-CAA-CBA

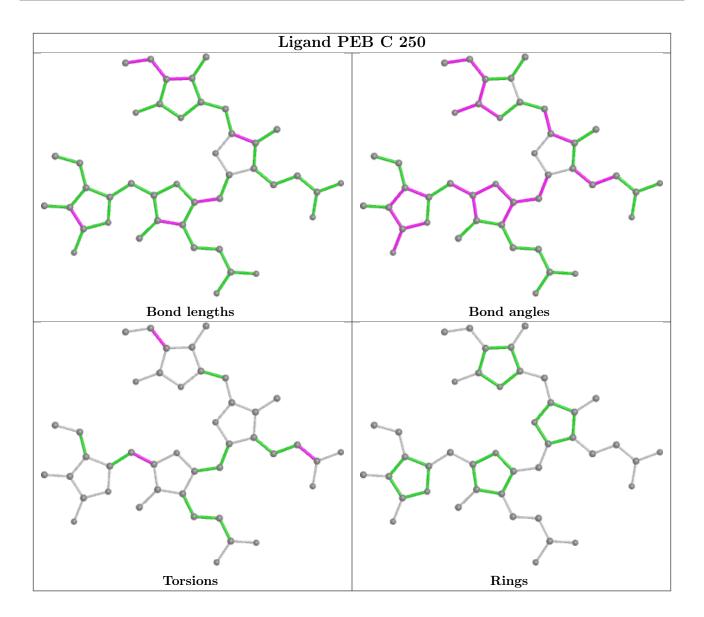
There are no ring outliers.

6 monomers are involved in 14 short contacts:

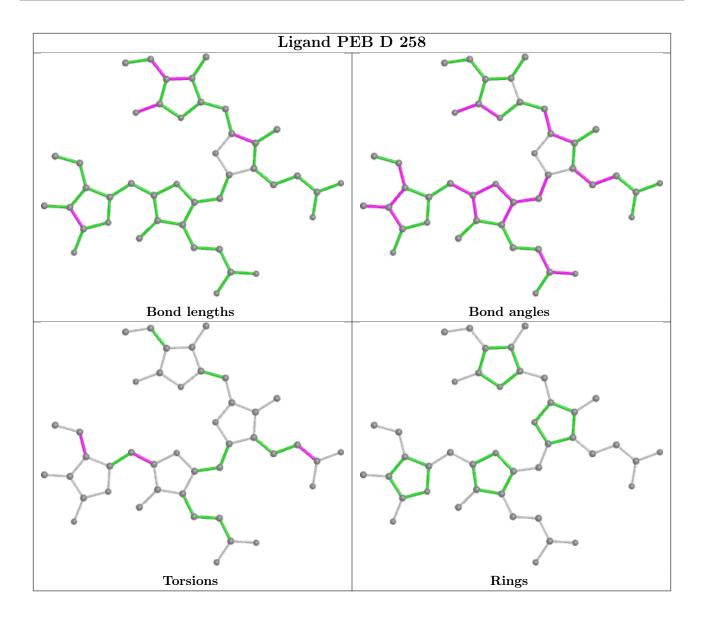
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	D	258	PEB	2	0
7	D	282	PEB	1	0
7	С	258	PEB	3	0
5	В	219	DBV	4	0
7	С	282	PEB	1	0
5	A	219	DBV	3	0

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 then 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.

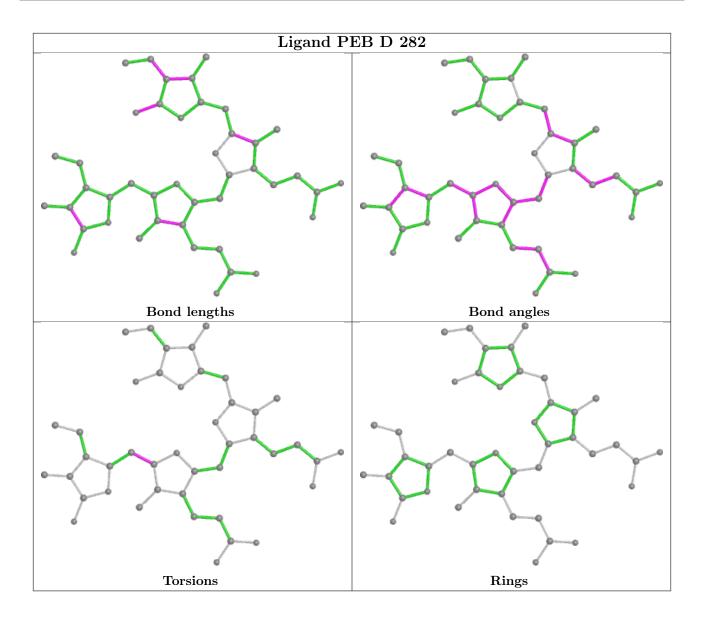




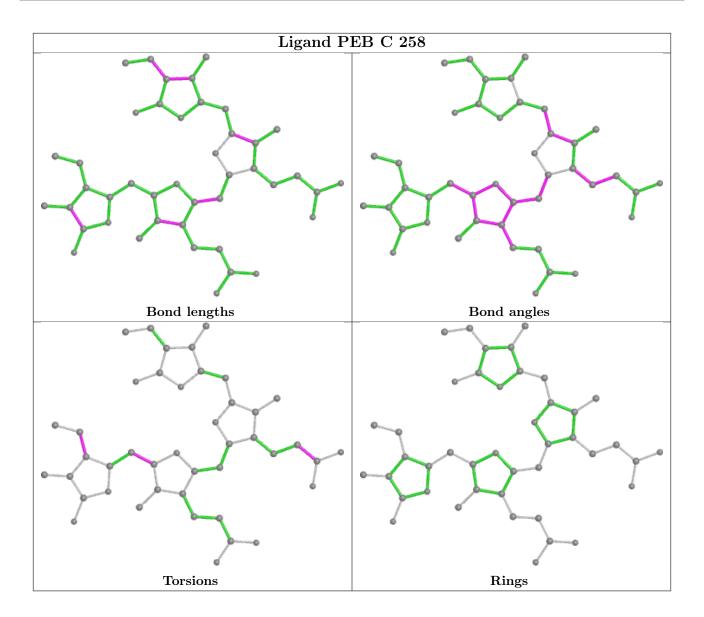




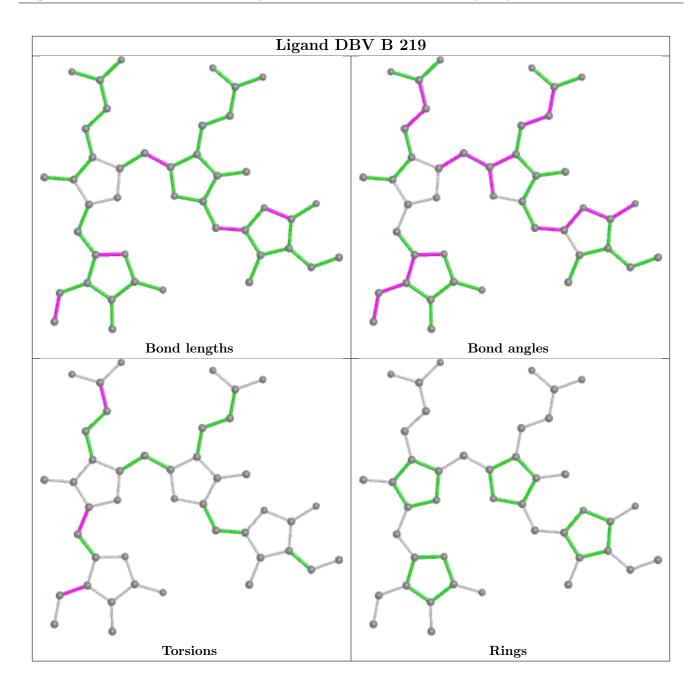




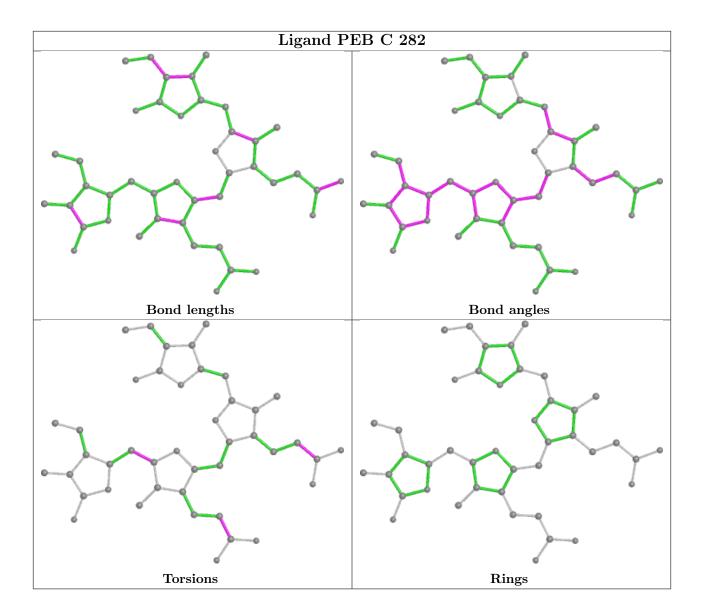




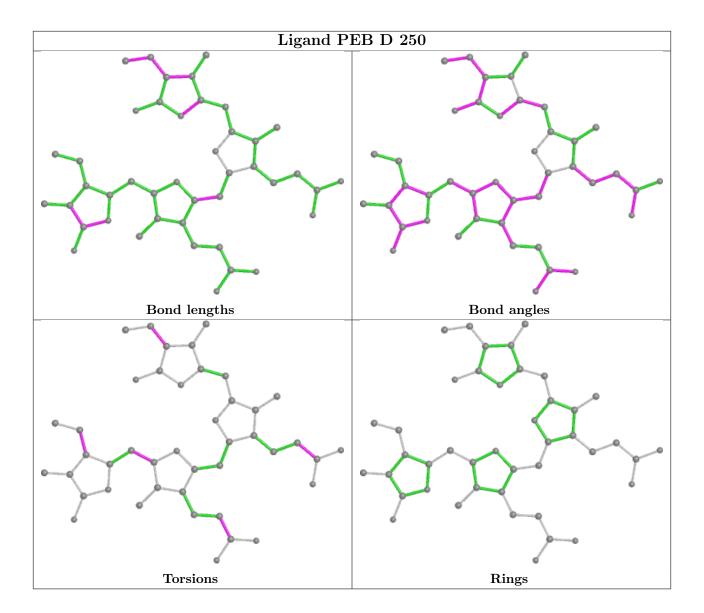




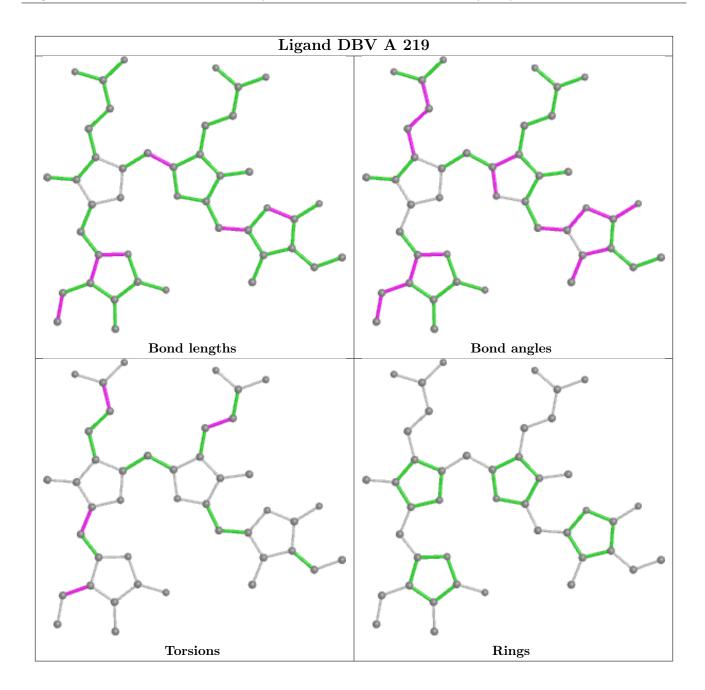












5.7 Other polymers (i)

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 (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.

