



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:38 PM UTC

PDB ID : 2C7J / pdb_00002c7j
Title : Phycoerythrocyanin from Mastigocladus laminosus, 295 K, 3.0 Å
Authors : Schmidt, M.; Krasselt, A.; Reuter, W.
Deposited on : 2005-11-24
Resolution : 3.00 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

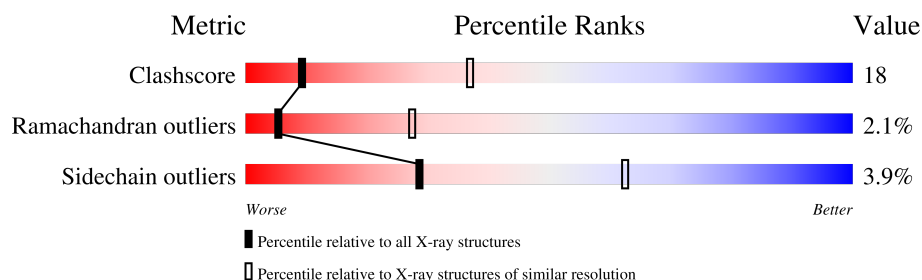
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)

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	162	
2	B	172	

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	CYC	A	1163	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2812 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.

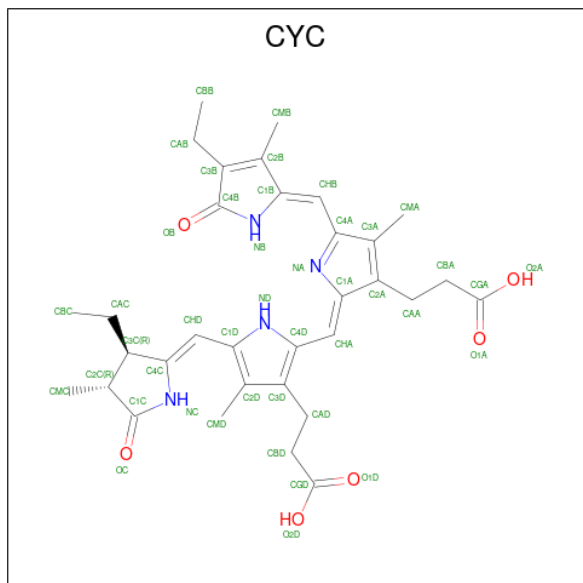
- Molecule 1 is a protein called Phycoerythrocyanin alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	162	Total	C	N	O	S	0	0	0
			1239	782	216	236	5			

- Molecule 2 is a protein called Phycoerythrocyanin beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	172	Total	C	N	O	S	0	0	0
			1295	801	234	252	8			

- Molecule 3 is PHYCOCYANOBILIN (CCD ID: CYC) (formula: $C_{33}H_{40}N_4O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			43	33	4	6		
3	B	1	Total	C	N	O	0	0
			43	33	4	6		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			43	33	4	6		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	68	Total	O	0	0
			68	68		
4	B	81	Total	O	0	0
			81	81		

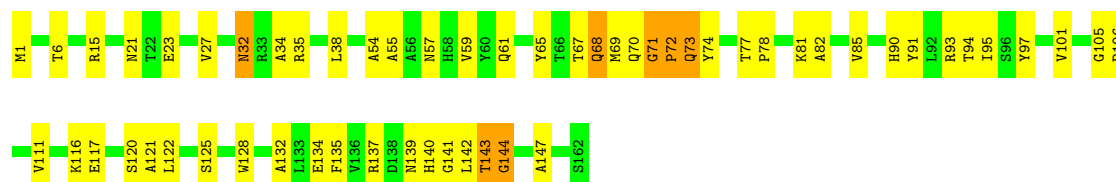
3 Residue-property plots

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.

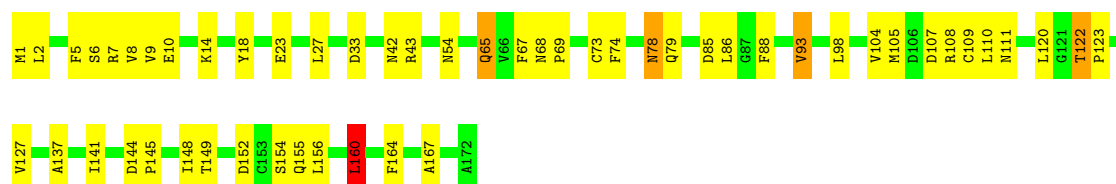
• Molecule 1: Phycoerythrocyanin alpha chain

Chain A: 



• Molecule 2: Phycoerythrocyanin beta chain

Chain B: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	156.75Å 156.75Å 40.20Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	51.30 – 3.00	Depositor
% Data completeness (in resolution range)	83.0 (51.30-3.00)	Depositor
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.189 , 0.260	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2812	wwPDB-VP
Average B, all atoms (Å ²)	33.0	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: CYC

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	0.45	0/1267	1.02	6/1720 (0.3%)
2	B	0.48	0/1312	1.05	5/1773 (0.3%)
All	All	0.46	0/2579	1.04	11/3493 (0.3%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	139	ASN	N-CA-C	8.48	120.52	111.28
1	A	71	GLY	N-CA-C	6.59	125.79	112.34
2	B	122	THR	CA-C-N	6.16	127.53	119.84
2	B	122	THR	C-N-CA	6.16	127.53	119.84
1	A	125	SER	CA-C-N	6.11	126.29	119.32
1	A	125	SER	C-N-CA	6.11	126.29	119.32
1	A	111	VAL	N-CA-C	5.97	119.38	113.71
2	B	109	CYS	N-CA-C	5.59	122.71	110.80
1	A	72	PRO	N-CA-C	-5.38	101.39	112.47
2	B	93	VAL	N-CA-C	-5.16	105.51	110.72
2	B	160	LEU	N-CA-C	-5.13	105.14	111.40

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	1239	0	1199	43	0
2	B	1295	0	1288	37	0
3	A	43	0	35	8	0
3	B	86	0	74	19	0
4	A	68	0	0	5	0
4	B	81	0	0	2	0
All	All	2812	0	2596	95	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 18.

All (95) 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)
3:B:1173:CYC:HMD3	3:B:1173:CYC:HC	1.35	0.92
3:B:1173:CYC:HC	3:B:1173:CYC:CMD	1.86	0.87
1:A:68:GLN:HG3	1:A:69:MET:H	1.48	0.79
1:A:73:GLN:HA	3:A:1163:CYC:HBD1	1.65	0.79
2:B:137:ALA:O	2:B:141:ILE:HG12	1.81	0.79
3:B:1173:CYC:HMD3	3:B:1173:CYC:NC	2.00	0.76
3:B:1174:CYC:CMD	3:B:1174:CYC:HC	1.99	0.76
2:B:1:MET:HG3	2:B:104:VAL:HG22	1.67	0.74
2:B:149:THR:O	3:B:1174:CYC:HMD2	1.87	0.74
1:A:143:THR:HG23	1:A:144:GLY:H	1.57	0.70
3:A:1163:CYC:HC	3:A:1163:CYC:HMD3	1.56	0.70
3:B:1174:CYC:HC	3:B:1174:CYC:HMD3	1.58	0.69
1:A:71:GLY:HA3	1:A:74:TYR:HB2	1.75	0.68
1:A:67:THR:HG23	1:A:81:LYS:HZ2	1.59	0.67
1:A:67:THR:HG23	1:A:81:LYS:NZ	2.09	0.67
1:A:68:GLN:HG3	1:A:69:MET:N	2.10	0.66
3:B:1174:CYC:HMB3	3:B:1174:CYC:HMA2	1.78	0.65
3:A:1163:CYC:HMD3	3:A:1163:CYC:NC	2.12	0.65
1:A:23:GLU:O	1:A:27:VAL:HG23	1.99	0.62
3:B:1174:CYC:HMD3	3:B:1174:CYC:NC	2.13	0.62
2:B:79:GLN:HE21	2:B:79:GLN:HA	1.66	0.61
1:A:73:GLN:HA	3:A:1163:CYC:CBD	2.31	0.61
3:A:1163:CYC:HC	3:A:1163:CYC:CMD	2.13	0.61
3:A:1163:CYC:HBC2	3:A:1163:CYC:CMC	2.32	0.60
1:A:90:HIS:CD2	1:A:93:ARG:HH21	2.20	0.59
1:A:72:PRO:HG2	1:A:73:GLN:HG3	1.83	0.58
1:A:21:ASN:HB2	4:A:2017:HOH:O	2.02	0.58
3:B:1174:CYC:HMA2	3:B:1174:CYC:CMB	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:ASN:N	1:A:32:ASN:HD22	2.02	0.58
1:A:142:LEU:HB3	1:A:147:ALA:HB2	1.86	0.57
1:A:91:TYR:O	1:A:95:ILE:HG13	2.05	0.56
3:B:1174:CYC:HMB3	3:B:1174:CYC:CMA	2.34	0.56
1:A:93:ARG:HA	2:B:18:TYR:CE2	2.41	0.55
1:A:15:ARG:HD3	4:A:2008:HOH:O	2.06	0.55
1:A:59:VAL:HB	1:A:132:ALA:HB2	1.89	0.55
2:B:155:GLN:HB3	4:B:2078:HOH:O	2.06	0.55
3:B:1174:CYC:HMB3	3:B:1174:CYC:C3A	2.37	0.54
3:A:1163:CYC:HBC2	3:A:1163:CYC:HMC3	1.90	0.54
2:B:105:MET:HG3	2:B:110:LEU:HD13	1.90	0.53
1:A:134:GLU:O	1:A:137:ARG:HB3	2.08	0.53
2:B:2:LEU:HD23	2:B:7:ARG:HE	1.73	0.53
2:B:5:PHE:O	2:B:8:VAL:HG12	2.10	0.52
3:B:1173:CYC:CMD	3:B:1173:CYC:NC	2.63	0.51
1:A:35:ARG:HB2	1:A:35:ARG:NH1	2.25	0.51
1:A:77:THR:HB	1:A:78:PRO:HD2	1.92	0.51
1:A:116:LYS:HD2	4:A:2055:HOH:O	2.11	0.51
3:B:1174:CYC:CMD	3:B:1174:CYC:NC	2.71	0.50
1:A:35:ARG:HB2	1:A:35:ARG:HH11	1.77	0.50
1:A:71:GLY:CA	1:A:74:TYR:HB2	2.41	0.50
3:B:1174:CYC:HC	3:B:1174:CYC:HMD2	1.77	0.49
1:A:34:ALA:O	1:A:38:LEU:HG	2.13	0.49
2:B:23:GLU:O	2:B:27:LEU:HD13	2.13	0.49
2:B:10:GLU:O	2:B:14:LYS:HG2	2.13	0.48
1:A:120:SER:HB3	4:A:2057:HOH:O	2.13	0.48
2:B:107:ASP:O	2:B:111:ASN:HB2	2.13	0.48
2:B:68:ASN:HB3	2:B:69:PRO:HD2	1.95	0.47
2:B:1:MET:HE2	2:B:104:VAL:HG13	1.96	0.47
2:B:110:LEU:HD21	2:B:167:ALA:HA	1.97	0.47
1:A:65:TYR:HB2	1:A:69:MET:HE3	1.96	0.47
2:B:54:ASN:N	2:B:54:ASN:HD22	2.13	0.47
2:B:152:ASP:OD2	2:B:154:SER:HB2	2.15	0.47
2:B:93:VAL:HG23	2:B:105:MET:HE1	1.96	0.46
1:A:69:MET:HE3	1:A:69:MET:HB2	1.77	0.46
2:B:86:LEU:HG	3:B:1173:CYC:HBC1	1.98	0.46
2:B:33:ASP:HB2	4:B:2029:HOH:O	2.16	0.45
2:B:68:ASN:HB3	2:B:69:PRO:CD	2.47	0.45
1:A:27:VAL:HG11	2:B:98:LEU:HD11	1.98	0.44
1:A:67:THR:HG22	1:A:67:THR:O	2.15	0.44
1:A:117:GLU:HB2	4:A:2056:HOH:O	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:LEU:HD12	1:A:122:LEU:HA	1.85	0.44
2:B:79:GLN:HA	2:B:79:GLN:NE2	2.32	0.44
2:B:67:PHE:HD2	2:B:79:GLN:HE22	1.66	0.44
3:A:1163:CYC:CMA	3:A:1163:CYC:NB	2.82	0.43
1:A:1:MET:HG2	2:B:6:SER:OG	2.18	0.43
1:A:55:ALA:O	1:A:59:VAL:HG12	2.18	0.43
1:A:101:VAL:HG23	2:B:9:VAL:HG21	2.01	0.43
2:B:108:ARG:O	3:B:1173:CYC:HBB1	2.19	0.43
1:A:82:ALA:O	1:A:85:VAL:HG12	2.19	0.42
1:A:71:GLY:HA2	1:A:72:PRO:HD2	1.88	0.42
1:A:140:HIS:ND1	1:A:141:GLY:N	2.67	0.42
2:B:122:THR:HA	2:B:123:PRO:HD3	1.91	0.42
2:B:160:LEU:HD22	2:B:164:PHE:CE2	2.54	0.42
1:A:94:THR:O	1:A:97:TYR:HB2	2.20	0.42
1:A:77:THR:O	1:A:81:LYS:HG3	2.20	0.42
2:B:85:ASP:O	2:B:88:PHE:HB2	2.19	0.42
1:A:54:ALA:HB1	1:A:135:PHE:CE1	2.54	0.41
2:B:74:PHE:O	2:B:78:ASN:HB2	2.20	0.41
2:B:43:ARG:NH2	2:B:148:ILE:HB	2.35	0.41
2:B:127:VAL:HG22	3:B:1173:CYC:H3C	2.02	0.41
2:B:65:GLN:CD	2:B:65:GLN:H	2.29	0.41
2:B:144:ASP:HA	2:B:145:PRO:HD3	1.90	0.41
1:A:128:TRP:CD1	1:A:128:TRP:N	2.88	0.40
2:B:148:ILE:HG21	3:B:1174:CYC:HMC3	2.03	0.40
1:A:57:ASN:HD22	1:A:57:ASN:N	2.20	0.40
2:B:73:CYS:HB2	3:B:1173:CYC:OC	2.21	0.40

There are no symmetry-related clashes.

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	160/162 (99%)	144 (90%)	10 (6%)	6 (4%)	2	15
2	B	170/172 (99%)	159 (94%)	10 (6%)	1 (1%)	21	56
All	All	330/334 (99%)	303 (92%)	20 (6%)	7 (2%)	5	27

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	68	GLN
1	A	143	THR
1	A	144	GLY
1	A	121	ALA
2	B	120	LEU
1	A	70	GLN
1	A	105	GLY

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	123/123 (100%)	118 (96%)	5 (4%)	27	61
2	B	135/135 (100%)	130 (96%)	5 (4%)	30	64
All	All	258/258 (100%)	248 (96%)	10 (4%)	28	62

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	A	32	ASN
1	A	61	GLN
1	A	73	GLN
1	A	106	PRO
2	B	42	ASN
2	B	65	GLN
2	B	78	ASN
2	B	156	LEU

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Mol	Chain	Res	Type
2	B	160	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	57	ASN
1	A	70	GLN
1	A	90	HIS
2	B	21	ASN
2	B	28	GLN
2	B	35	ASN
2	B	54	ASN
2	B	65	GLN
2	B	78	ASN
2	B	79	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](#)

3 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	CYC	B	1173	2	46,46,46	2.30	12 (26%)	63,67,67	2.38	18 (28%)
3	CYC	A	1163	1	46,46,46	2.50	11 (23%)	63,67,67	2.79	17 (26%)
3	CYC	B	1174	2	46,46,46	2.19	12 (26%)	63,67,67	2.34	15 (23%)

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	CYC	B	1173	2	-	10/26/74/74	0/4/4/4
3	CYC	A	1163	1	2/2/14/19	15/26/74/74	0/4/4/4
3	CYC	B	1174	2	-	14/26/74/74	0/4/4/4

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1163	CYC	CHD-C4C	10.34	1.53	1.36
3	B	1173	CYC	CHD-C4C	10.00	1.53	1.36
3	B	1174	CYC	CHD-C4C	9.21	1.51	1.36
3	A	1163	CYC	C4C-NC	6.94	1.50	1.37
3	A	1163	CYC	CHD-C1D	5.02	1.51	1.40
3	A	1163	CYC	C2C-C3C	-4.70	1.41	1.54
3	B	1173	CYC	CHD-C1D	4.51	1.50	1.40
3	B	1173	CYC	C4C-NC	4.22	1.45	1.37
3	B	1174	CYC	C4C-NC	4.14	1.45	1.37
3	B	1173	CYC	C2C-C3C	-4.06	1.43	1.54
3	B	1173	CYC	C1C-NC	-3.99	1.32	1.37
3	B	1174	CYC	C2C-C3C	-3.45	1.45	1.54
3	B	1174	CYC	CHD-C1D	3.28	1.47	1.40
3	B	1174	CYC	C1C-NC	-3.17	1.33	1.37
3	A	1163	CYC	C1B-C2B	-2.89	1.39	1.45
3	B	1174	CYC	CHA-C1A	2.80	1.44	1.38
3	A	1163	CYC	C2C-C1C	-2.72	1.49	1.52
3	A	1163	CYC	CBD-CGD	2.56	1.56	1.50
3	B	1174	CYC	CMD-C2D	-2.53	1.45	1.50
3	A	1163	CYC	O1D-CGD	2.44	1.30	1.22
3	B	1174	CYC	C3C-C4C	-2.37	1.45	1.50
3	B	1174	CYC	C1B-NB	2.29	1.41	1.37
3	B	1173	CYC	O1D-CGD	2.29	1.29	1.22
3	B	1173	CYC	C1B-NB	2.28	1.41	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1173	CYC	CHA-C1A	2.28	1.43	1.38
3	B	1174	CYC	CBD-CGD	2.28	1.55	1.50
3	A	1163	CYC	CHA-C1A	2.24	1.42	1.38
3	B	1174	CYC	C4B-NB	-2.20	1.32	1.38
3	A	1163	CYC	CMB-C2B	-2.16	1.46	1.50
3	B	1173	CYC	C4B-NB	-2.14	1.33	1.38
3	B	1173	CYC	CMD-C2D	-2.10	1.46	1.50
3	B	1174	CYC	C4A-NA	2.09	1.41	1.36
3	A	1163	CYC	C4B-NB	-2.08	1.33	1.38
3	B	1173	CYC	CBD-CGD	2.06	1.55	1.50
3	B	1173	CYC	C1B-C2B	-2.00	1.41	1.45

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1163	CYC	CHD-C4C-NC	-12.23	110.37	125.63
3	B	1174	CYC	C1C-NC-C4C	-8.32	102.97	113.41
3	B	1173	CYC	C1C-NC-C4C	-8.00	103.38	113.41
3	B	1173	CYC	C2C-C1C-NC	7.80	114.78	108.29
3	B	1174	CYC	C2C-C1C-NC	7.66	114.66	108.29
3	A	1163	CYC	CMC-C2C-C1C	6.41	126.22	112.40
3	A	1163	CYC	C1D-CHD-C4C	-6.02	117.49	127.76
3	B	1173	CYC	C3B-C4B-NB	5.71	111.32	106.77
3	A	1163	CYC	C3B-C4B-NB	5.67	111.29	106.77
3	B	1174	CYC	C3B-C4B-NB	5.52	111.17	106.77
3	A	1163	CYC	C2C-C3C-C4C	5.47	109.53	101.34
3	A	1163	CYC	CAC-C3C-C4C	5.29	126.26	112.67
3	A	1163	CYC	C1B-NB-C4B	-5.04	104.47	110.66
3	B	1173	CYC	C1B-NB-C4B	-4.77	104.80	110.66
3	B	1174	CYC	C3C-C4C-NC	4.62	113.89	107.94
3	B	1173	CYC	CMD-C2D-C1D	4.30	131.87	125.62
3	B	1174	CYC	OC-C1C-C2C	-4.15	122.87	126.17
3	B	1174	CYC	C1B-NB-C4B	-3.99	105.76	110.66
3	B	1173	CYC	C3C-C4C-NC	3.92	112.99	107.94
3	B	1174	CYC	CMD-C2D-C1D	3.87	131.24	125.62
3	A	1163	CYC	OB-C4B-C3B	-3.77	124.07	128.03
3	A	1163	CYC	CAB-C3B-C4B	3.62	126.97	121.37
3	B	1173	CYC	OB-C4B-C3B	-3.62	124.23	128.03
3	B	1173	CYC	OC-C1C-C2C	-3.58	123.32	126.17
3	B	1174	CYC	CAB-C3B-C4B	3.57	126.89	121.37
3	A	1163	CYC	C2B-C1B-NB	3.47	112.03	106.97
3	A	1163	CYC	CMC-C2C-C3C	3.47	127.39	113.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1173	CYC	CAB-C3B-C4B	3.45	126.70	121.37
3	B	1174	CYC	OB-C4B-C3B	-3.41	124.44	128.03
3	A	1163	CYC	CMD-C2D-C1D	3.41	130.58	125.62
3	B	1173	CYC	C2B-C1B-NB	3.33	111.82	106.97
3	B	1174	CYC	CAD-C3D-C4D	3.20	130.84	125.77
3	A	1163	CYC	C3C-C4C-NC	-3.14	103.90	107.94
3	B	1173	CYC	CAD-C3D-C4D	3.02	130.57	125.77
3	A	1163	CYC	C1C-NC-C4C	-2.89	109.78	113.41
3	B	1174	CYC	C2B-C1B-NB	2.76	110.99	106.97
3	B	1173	CYC	CBC-CAC-C3C	-2.73	107.62	113.41
3	B	1174	CYC	CMB-C2B-C1B	2.68	127.42	124.16
3	B	1174	CYC	CAC-C3C-C4C	-2.61	105.98	112.67
3	B	1173	CYC	OC-C1C-NC	-2.57	121.90	124.93
3	A	1163	CYC	C2C-C1C-NC	2.54	110.40	108.29
3	A	1163	CYC	CAD-C3D-C4D	2.46	129.67	125.77
3	A	1163	CYC	CHB-C1B-C2B	-2.43	122.13	126.97
3	B	1173	CYC	CMB-C2B-C1B	2.30	126.95	124.16
3	B	1173	CYC	C4D-CHA-C1A	-2.30	123.21	128.22
3	B	1173	CYC	C1B-CHB-C4A	-2.24	122.56	128.06
3	B	1174	CYC	CAD-C3D-C2D	-2.21	122.11	127.07
3	B	1173	CYC	CAD-C3D-C2D	-2.20	122.14	127.07
3	B	1173	CYC	CAD-CBD-CGD	-2.10	108.10	113.67
3	B	1174	CYC	OC-C1C-NC	-2.09	122.47	124.93

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	1163	CYC	C2C
3	A	1163	CYC	C3C

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1163	CYC	C4C-C3C-CAC-CBC
3	A	1163	CYC	NC-C4C-CHD-C1D
3	B	1173	CYC	NA-C4A-CHB-C1B
3	B	1173	CYC	C3A-C4A-CHB-C1B
3	B	1174	CYC	NB-C1B-CHB-C4A
3	B	1174	CYC	C2B-C1B-CHB-C4A
3	B	1174	CYC	C2C-C3C-CAC-CBC
3	B	1174	CYC	C4C-C3C-CAC-CBC
3	A	1163	CYC	ND-C1D-CHD-C4C

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Mol	Chain	Res	Type	Atoms
3	B	1173	CYC	C2D-C1D-CHD-C4C
3	B	1174	CYC	C2D-C1D-CHD-C4C
3	B	1173	CYC	ND-C1D-CHD-C4C
3	B	1174	CYC	ND-C1D-CHD-C4C
3	A	1163	CYC	C2D-C1D-CHD-C4C
3	A	1163	CYC	NA-C4A-CHB-C1B
3	B	1174	CYC	NA-C4A-CHB-C1B
3	A	1163	CYC	NB-C1B-CHB-C4A
3	A	1163	CYC	C3A-C4A-CHB-C1B
3	B	1174	CYC	C3A-C4A-CHB-C1B
3	A	1163	CYC	C2B-C1B-CHB-C4A
3	A	1163	CYC	C3D-CAD-CBD-CGD
3	B	1173	CYC	NA-C1A-CHA-C4D
3	B	1173	CYC	C2A-C1A-CHA-C4D
3	A	1163	CYC	NA-C1A-CHA-C4D
3	B	1174	CYC	CAD-CBD-CGD-O2D
3	B	1173	CYC	CAD-CBD-CGD-O2D
3	B	1174	CYC	CAD-CBD-CGD-O1D
3	B	1173	CYC	CAD-CBD-CGD-O1D
3	A	1163	CYC	CAA-CBA-CGA-O1A
3	A	1163	CYC	CAA-CBA-CGA-O2A
3	B	1174	CYC	NA-C1A-CHA-C4D
3	A	1163	CYC	CAD-CBD-CGD-O2D
3	B	1173	CYC	CAA-CBA-CGA-O2A
3	B	1174	CYC	CAA-CBA-CGA-O2A
3	B	1174	CYC	CAA-CBA-CGA-O1A
3	A	1163	CYC	C2A-C1A-CHA-C4D
3	B	1173	CYC	CAA-CBA-CGA-O1A
3	A	1163	CYC	CAD-CBD-CGD-O1D
3	B	1174	CYC	C2B-C3B-CAB-CBB

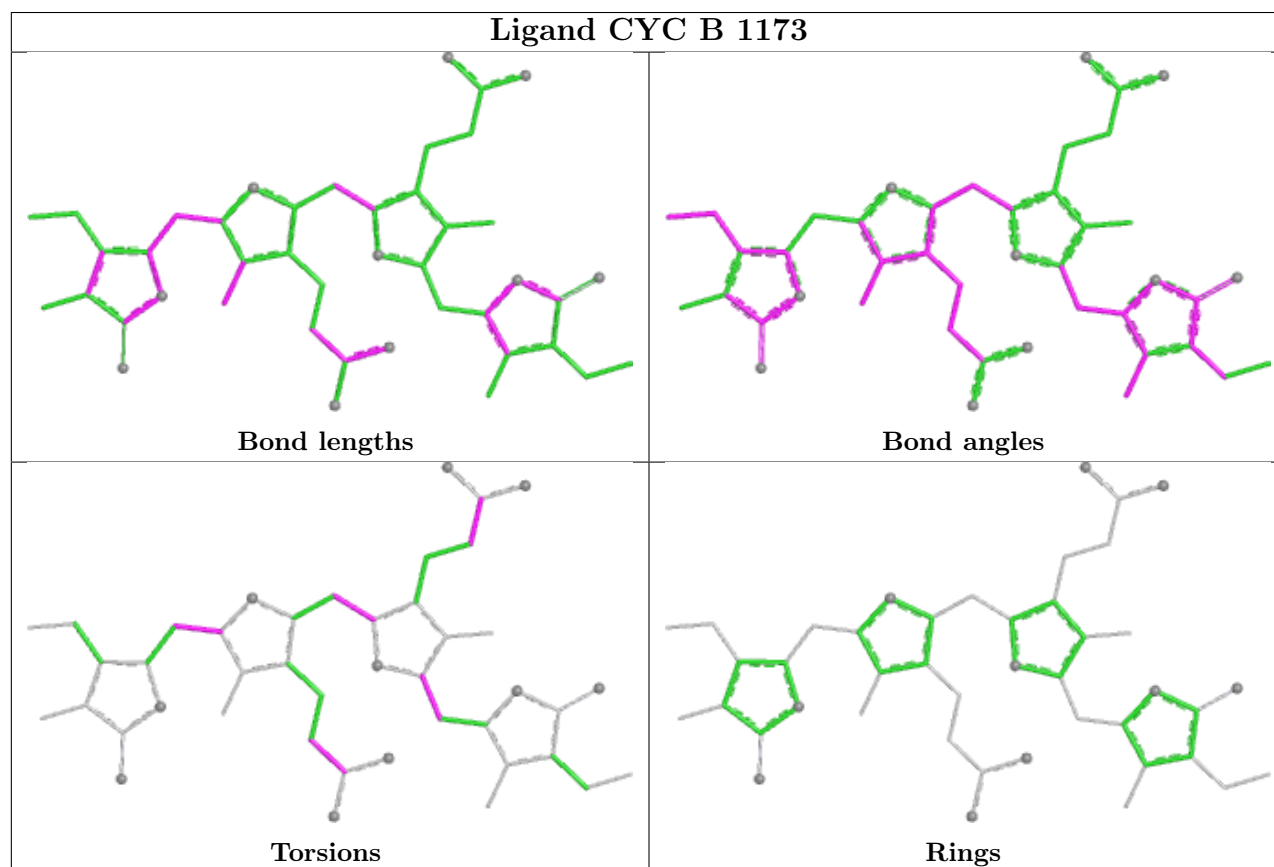
There are no ring outliers.

3 monomers are involved in 27 short contacts:

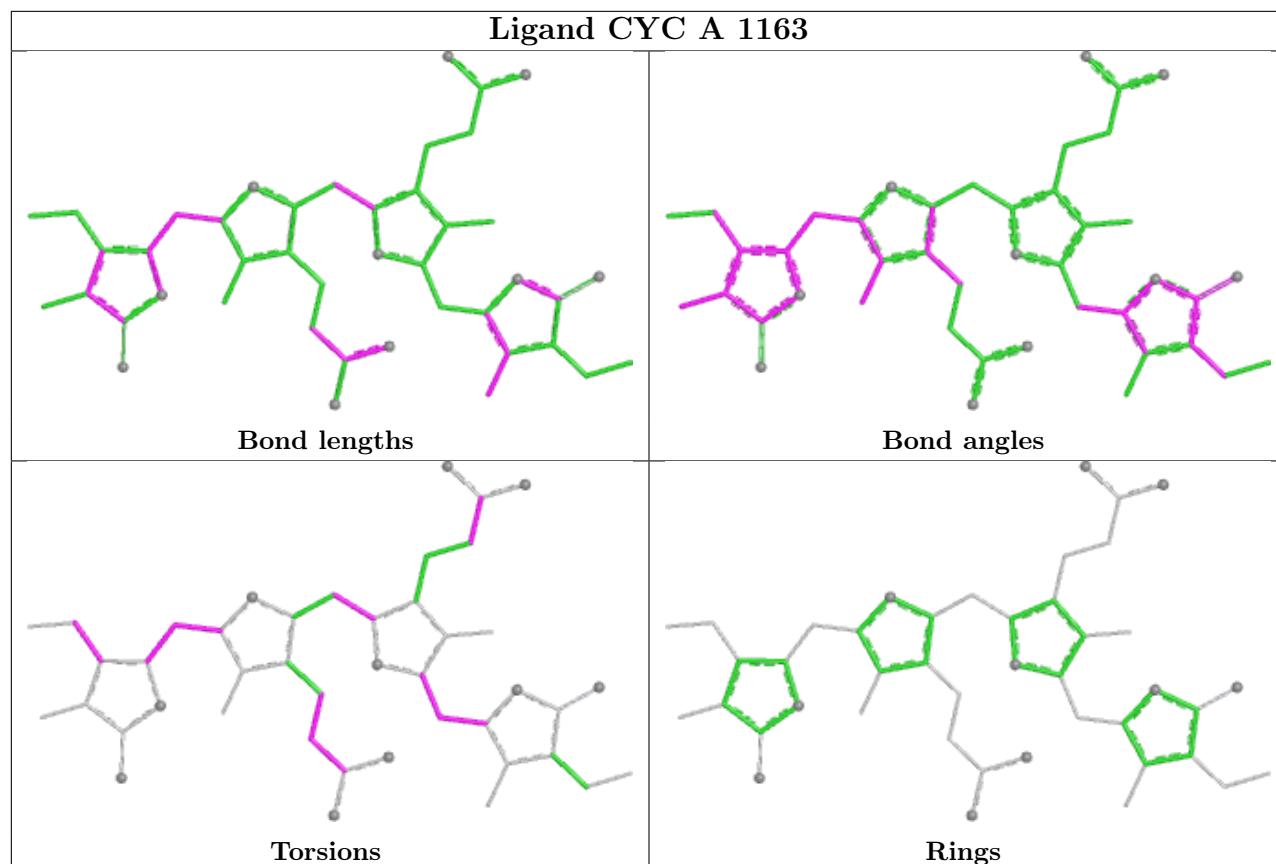
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1173	CYC	8	0
3	A	1163	CYC	8	0
3	B	1174	CYC	11	0

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

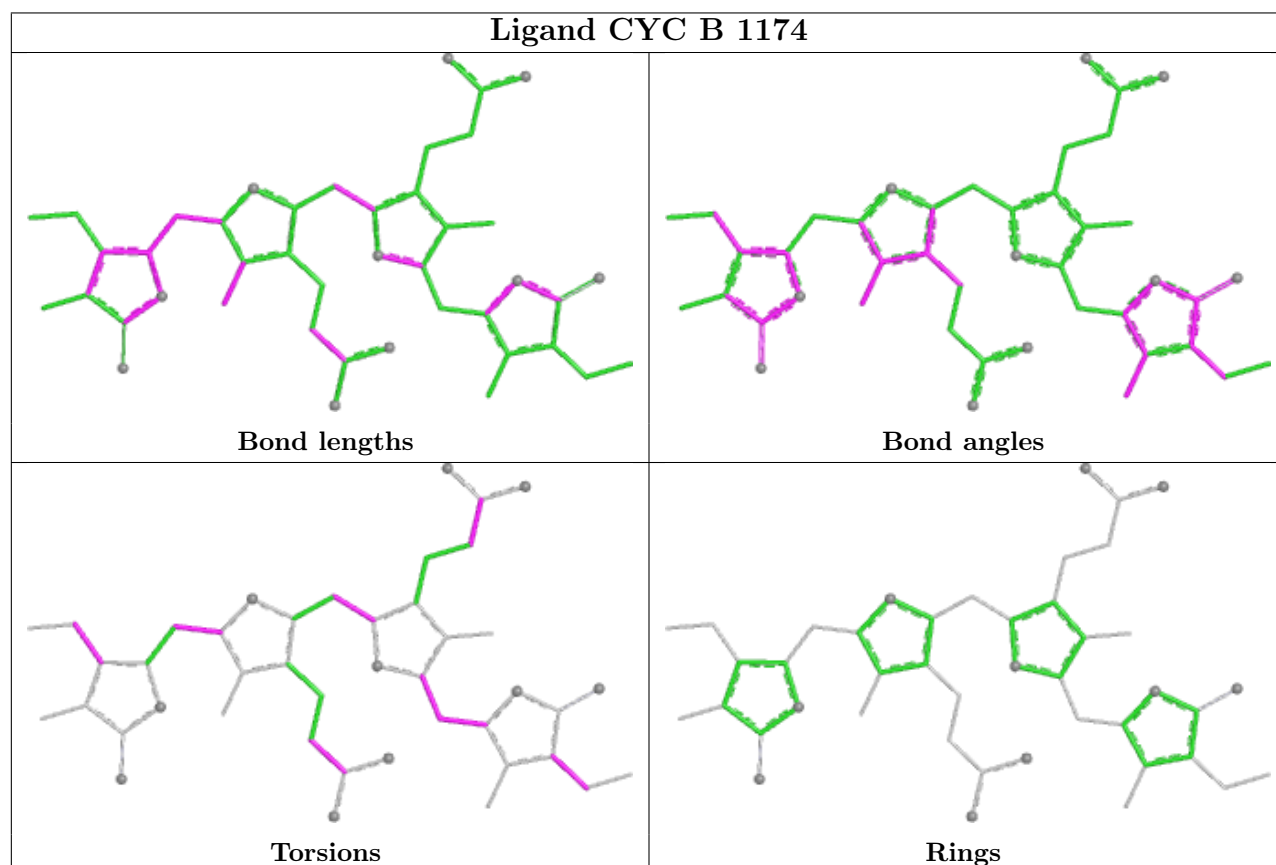
also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted 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 CYC A 1163



Ligand CYC B 1174



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.