



wwPDB X-ray Structure Validation Summary Report

Feb 26, 2014 – 07:07 PM GMT

PDB ID : 1FE1
Title : CRYSTAL STRUCTURE PHOTOSYSTEM II
Authors : Zouni, A.; Witt, H.-T.; Kern, J.; Fromme, P.; Krauss, N.; Saenger, W.; Orth, P.
Deposited on : 2000-07-20
Resolution : 3.80 Å(reported)

This is a wwPDB validation summary report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

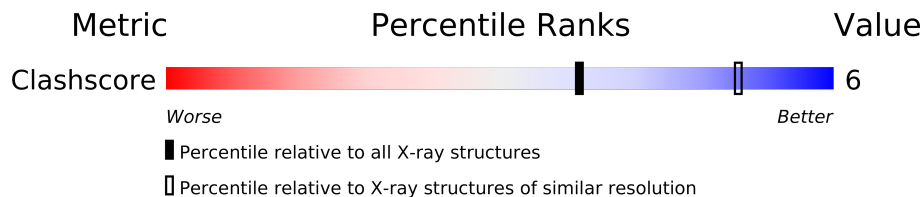
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	21963
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.80 Å.






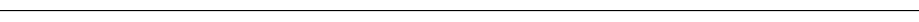








Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	79885	1100 (4.10-3.50)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	169	
1	J	169	
2	B	174	
2	K	174	
3	C	156	
3	L	156	
4	D	155	
4	M	155	
5	E	40	
5	N	40	
6	F	30	
6	O	30	
7	G	312	
7	P	312	
8	H	115	
8	Q	115	
9	I	87	
9	R	87	

2 Entry composition

There are 17 unique types of molecules in this entry. The entry contains 4328 atoms, of which 0 are hydrogen and 0 are deuterium.

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 (PHOTOSYSTEM II: SUBUNIT PSBA).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace
1	A	169	Total	C	0	0	169
			169	169			
1	J	169	Total	C	0	0	169
			169	169			

- Molecule 2 is a protein called PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBD).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace
2	B	174	Total	C	0	0	174
			174	174			
2	K	174	Total	C	0	0	174
			174	174			

- Molecule 3 is a protein called PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBC).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace
3	C	156	Total	C	0	0	156
			156	156			
3	L	156	Total	C	0	0	156
			156	156			

- Molecule 4 is a protein called PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBB).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace
4	D	155	Total	C	0	0	155
			155	155			
4	M	155	Total	C	0	0	155
			155	155			

- Molecule 5 is a protein called PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBE).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
5	E	40	Total C 40 40	0	0	40
5	N	40	Total C 40 40	0	0	40

- Molecule 6 is a protein called PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBF).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
6	F	30	Total C 30 30	0	0	30
6	O	30	Total C 30 30	0	0	30

- Molecule 7 is a protein called PROTEIN (PHOTOSYSTEM II: SUBUNIT UNKNOWN).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
7	G	312	Total C 312 312	0	0	312
7	P	312	Total C 312 312	0	0	312

- Molecule 8 is a protein called PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBO).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
8	H	115	Total C 115 115	0	0	115
8	Q	115	Total C 115 115	0	0	115

- Molecule 9 is a protein called PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBV).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
9	I	87	Total C 87 87	0	0	87
9	R	87	Total C 87 87	0	0	87

- Molecule 10 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	J	4	Total Mn 4 4	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	4	Total	Mn	0	0
			4	4		

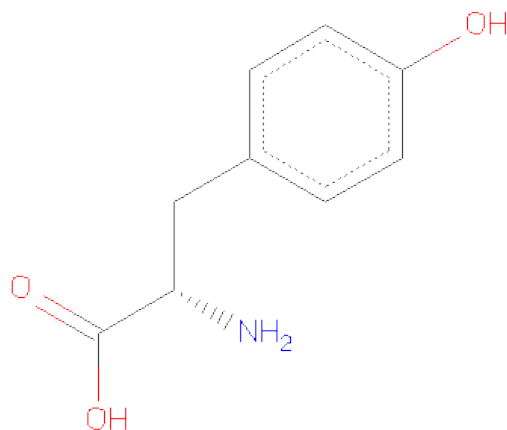
- Molecule 11 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	B	1	Total	Fe	0	0
			1	1		
11	K	1	Total	Fe	0	0
			1	1		

- Molecule 12 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	H	1	Total	Cd	0	0
			1	1		
12	Q	1	Total	Cd	0	0
			1	1		

- Molecule 13 is TYROSINE (three-letter code: TYR) (formula: C₉H₁₁NO₃).



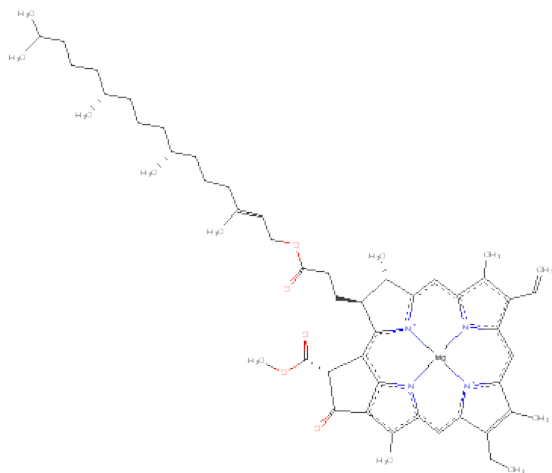
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
13	A	1	Total	C	O	0	0
			8	7	1		
13	B	1	Total	C	O	0	0
			8	7	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
13	J	1	Total	C	O	0	0
			8	7	1		
13	K	1	Total	C	O	0	0
			8	7	1		

- Molecule 14 is CHLOROPHYLL A (three-letter code: CLA) (formula: C₅₅H₇₂MgN₄O₅).



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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
14	C	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	C	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	C	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	C	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	C	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	C	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	C	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	C	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	G	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	D	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	D	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	D	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	D	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	D	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	D	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	D	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	D	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	D	1	Total	C	Mg	N	0	0
			25	20	1	4		

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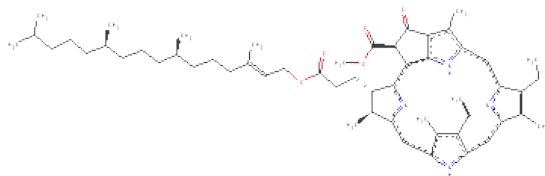
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
14	D	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	D	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	J	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	K	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	J	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	J	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	J	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	K	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	L	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	L	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	L	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	L	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	L	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	L	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	L	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	L	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	L	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	L	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	P	1	Total	C	Mg	N	0	0
			25	20	1	4		

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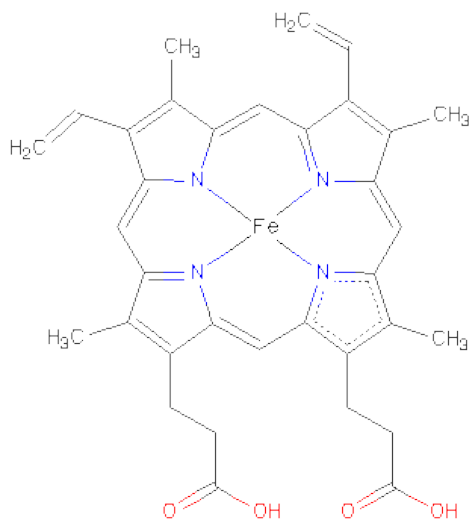
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
14	M	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	M	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	M	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	M	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	M	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	M	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	M	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	M	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	M	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	M	1	Total	C	Mg	N	0	0
			25	20	1	4		
14	M	1	Total	C	Mg	N	0	0
			25	20	1	4		

- Molecule 15 is PHEOPHYTIN A (three-letter code: PHO) (formula: C₅₅H₇₄N₄O₅).



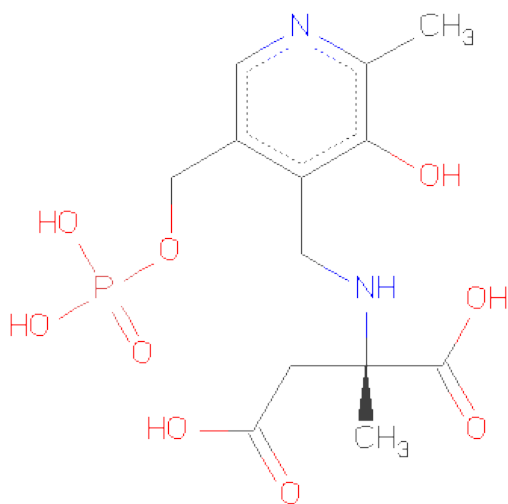
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
15	B	1	Total	C	N	0	0
			24	20	4		
15	A	1	Total	C	N	0	0
			24	20	4		
15	K	1	Total	C	N	0	0
			24	20	4		
15	J	1	Total	C	N	0	0
			24	20	4		

- Molecule 16 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
16	F	1	Total	C	Fe	N	0	0
			25	20	1	4		
16	I	1	Total	C	Fe	N	0	0
			25	20	1	4		
16	O	1	Total	C	Fe	N	0	0
			25	20	1	4		
16	R	1	Total	C	Fe	N	0	0
			25	20	1	4		

- Molecule 17 is 2-[(3-HYDROXY-2-METHYL-5-PHOSPHONOOXYMETHYL-PYRIDIN-4-YLMETHYL)-AMINO]-2-METHYL-SUCCINICACID (three-letter code: PLA) (formula: $C_{13}H_{19}N_2O_9P$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
17	B	1	Total	C	N	0	0
			6	5	1		
17	K	1	Total	C	N	0	0
			6	5	1		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

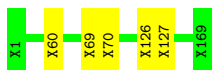
- Molecule 1: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBA)

Chain A: 



- Molecule 1: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBA)

Chain J: 



- Molecule 2: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBD)

Chain B: 



- Molecule 2: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBD)

Chain K: 



- Molecule 3: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBC)

Chain C: 



- Molecule 3: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBC)

Chain L: 



- Molecule 4: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBB)

Chain D: 

There are no outlier residues recorded for this chain.

- Molecule 4: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBB)

Chain M: 

There are no outlier residues recorded for this chain.

- Molecule 5: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBE)

Chain E: 

There are no outlier residues recorded for this chain.

- Molecule 5: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBE)

Chain N: 

There are no outlier residues recorded for this chain.

- Molecule 6: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBF)

Chain F: 

There are no outlier residues recorded for this chain.

- Molecule 6: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBF)

Chain O: 

There are no outlier residues recorded for this chain.

- Molecule 7: PROTEIN (PHOTOSYSTEM II: SUBUNIT UNKNOWN)

Chain G: 

There are no outlier residues recorded for this chain.

- Molecule 7: PROTEIN (PHOTOSYSTEM II: SUBUNIT UNKNOWN)

Chain P: 

There are no outlier residues recorded for this chain.

- Molecule 8: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBO)

Chain H: 

There are no outlier residues recorded for this chain.

- Molecule 8: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBO)

Chain Q: 

There are no outlier residues recorded for this chain.

- Molecule 9: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBV)

Chain I: 

There are no outlier residues recorded for this chain.

- Molecule 9: PROTEIN (PHOTOSYSTEM II: SUBUNIT PSBV)

Chain R: 

There are no outlier residues recorded for this chain.

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	130.01Å 226.72Å 308.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.80	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-3.80)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	?	Depositor
R, R_{free}	(Not available) , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4328	wwPDB-VP
Average B, all atoms (Å ²)	28.0	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: PHO, MN, CLA, CD, FE, PLA, HEM

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	$\# Z > 5$

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	169	0	0	4	0
1	J	169	0	0	4	0
2	B	174	0	0	4	0
2	K	174	0	0	4	0
3	C	156	0	0	2	0
3	L	156	0	0	2	0
4	D	155	0	0	0	0
4	M	155	0	0	0	0
5	E	40	0	0	0	0
5	N	40	0	0	0	0
6	F	30	0	0	0	0
6	O	30	0	0	0	0
7	G	312	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	P	312	0	0	0	0
8	H	115	0	0	0	0
8	Q	115	0	0	0	0
9	I	87	0	0	0	0
9	R	87	0	0	0	0
10	A	4	0	0	0	0
10	J	4	0	0	0	0
11	B	1	0	0	0	0
11	K	1	0	0	0	0
12	H	1	0	0	0	0
12	Q	1	0	0	0	0
13	A	8	0	5	0	0
13	B	8	0	5	0	0
13	J	8	0	5	0	0
13	K	8	0	5	0	0
14	A	100	0	12	0	0
14	B	50	0	6	0	0
14	C	300	0	36	2	0
14	D	325	0	39	6	0
14	G	25	0	3	0	0
14	J	100	0	12	0	0
14	K	50	0	6	0	0
14	L	300	0	36	2	0
14	M	325	0	39	6	0
14	P	25	0	3	0	0
15	A	24	0	5	0	0
15	B	24	0	5	0	0
15	J	24	0	5	0	0
15	K	24	0	5	0	0
16	F	25	0	4	0	0
16	I	25	0	4	0	0
16	O	25	0	4	0	0
16	R	25	0	4	0	0
17	B	6	0	1	0	0
17	K	6	0	1	0	0
All	All	4328	0	250	26	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

The worst 5 of 26 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:127:UNK:CA	2:K:53:UNK:CA	1.88	1.51
1:A:127:UNK:CA	2:B:53:UNK:CA	1.88	1.49
1:A:60:UNK:CA	2:B:131:UNK:CA	1.95	1.45
1:J:60:UNK:CA	2:K:131:UNK:CA	1.95	1.44
3:L:149:UNK:CA	14:L:166:CLA:C2A	2.23	1.17

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains ⓘ

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

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 90 ligands modelled in this entry, 12 are monoatomic - leaving 78 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	$\# Z > 2$
13	TYR	A	174	1	8,8,13	1.10	1 (12%)	10,10,17	1.33	2 (20%)
14	CLA	A	175	-	23,32,73	7.68	5 (21%)	12,54,113	1.71	2 (16%)
14	CLA	A	176	-	23,32,73	7.60	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	A	177	-	23,32,73	7.52	5 (21%)	12,54,113	1.70	2 (16%)
15	PHO	A	178	-	22,28,69	31.63	13 (59%)	8,40,99	1.44	1 (12%)
14	CLA	A	179	-	23,32,73	7.58	5 (21%)	12,54,113	1.70	2 (16%)
13	TYR	B	176	2	8,8,13	1.10	1 (12%)	10,10,17	1.33	2 (20%)
14	CLA	B	177	-	23,32,73	7.56	5 (21%)	12,54,113	1.69	2 (16%)
15	PHO	B	178	-	22,28,69	31.79	13 (59%)	8,40,99	1.42	1 (12%)
14	CLA	B	179	-	23,32,73	7.59	5 (21%)	12,54,113	1.71	2 (16%)
17	PLA	B	180	-	5,6,25	15.03	3 (60%)	1,6,37	0.65	0
14	CLA	C	157	-	23,32,73	7.45	5 (21%)	12,54,113	1.72	2 (16%)
14	CLA	C	158	-	23,32,73	7.59	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	C	159	-	23,32,73	7.64	5 (21%)	12,54,113	1.68	2 (16%)
14	CLA	C	160	-	23,32,73	7.56	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	C	161	-	23,32,73	7.51	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	C	162	-	23,32,73	7.56	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	C	163	-	23,32,73	7.61	5 (21%)	12,54,113	1.71	2 (16%)
14	CLA	C	164	-	23,32,73	7.61	5 (21%)	12,54,113	1.68	2 (16%)
14	CLA	C	165	-	23,32,73	7.62	5 (21%)	12,54,113	1.71	2 (16%)
14	CLA	C	166	-	23,32,73	7.38	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	C	167	-	23,32,73	7.58	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	C	168	-	23,32,73	7.52	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	D	156	-	23,32,73	7.52	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	D	157	-	23,32,73	7.57	5 (21%)	12,54,113	1.71	2 (16%)
14	CLA	D	158	-	23,32,73	7.63	5 (21%)	12,54,113	1.72	2 (16%)
14	CLA	D	159	-	23,32,73	7.52	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	D	160	-	23,32,73	7.61	5 (21%)	12,54,113	1.68	2 (16%)
14	CLA	D	161	-	23,32,73	7.56	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	D	162	-	23,32,73	7.58	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	D	163	-	23,32,73	7.54	5 (21%)	12,54,113	1.68	2 (16%)
14	CLA	D	164	-	23,32,73	7.47	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	D	165	-	23,32,73	7.68	5 (21%)	12,54,113	1.68	2 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
14	CLA	D	166	-	23,32,73	7.63	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	D	167	-	23,32,73	7.55	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	D	168	-	23,32,73	7.60	5 (21%)	12,54,113	1.70	2 (16%)
16	HEM	F	31	-	22,32,50	30.36	6 (27%)	16,54,82	1.51	3 (18%)
14	CLA	G	313	-	23,32,73	7.51	5 (21%)	12,54,113	1.70	2 (16%)
16	HEM	I	88	-	22,32,50	30.41	6 (27%)	16,54,82	1.53	3 (18%)
13	TYR	J	174	1	8,8,13	1.10	1 (12%)	10,10,17	1.32	2 (20%)
14	CLA	J	175	-	23,32,73	7.64	5 (21%)	12,54,113	1.71	2 (16%)
14	CLA	J	176	-	23,32,73	7.60	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	J	177	-	23,32,73	7.50	5 (21%)	12,54,113	1.70	2 (16%)
15	PHO	J	178	-	22,28,69	31.65	13 (59%)	8,40,99	1.44	1 (12%)
14	CLA	J	179	-	23,32,73	7.64	5 (21%)	12,54,113	1.69	2 (16%)
13	TYR	K	176	2	8,8,13	1.10	1 (12%)	10,10,17	1.33	2 (20%)
14	CLA	K	177	-	23,32,73	7.58	5 (21%)	12,54,113	1.69	2 (16%)
15	PHO	K	178	-	22,28,69	31.64	13 (59%)	8,40,99	1.41	1 (12%)
14	CLA	K	179	-	23,32,73	7.67	5 (21%)	12,54,113	1.70	2 (16%)
17	PLA	K	180	-	5,6,25	15.03	3 (60%)	1,6,37	0.65	0
14	CLA	L	157	-	23,32,73	7.48	5 (21%)	12,54,113	1.73	2 (16%)
14	CLA	L	158	-	23,32,73	7.57	5 (21%)	12,54,113	1.71	2 (16%)
14	CLA	L	159	-	23,32,73	7.65	5 (21%)	12,54,113	1.68	2 (16%)
14	CLA	L	160	-	23,32,73	7.52	5 (21%)	12,54,113	1.71	2 (16%)
14	CLA	L	161	-	23,32,73	7.46	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	L	162	-	23,32,73	7.46	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	L	163	-	23,32,73	7.66	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	L	164	-	23,32,73	7.61	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	L	165	-	23,32,73	7.69	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	L	166	-	23,32,73	7.40	5 (21%)	12,54,113	1.71	2 (16%)
14	CLA	L	167	-	23,32,73	7.54	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	L	168	-	23,32,73	7.46	5 (21%)	12,54,113	1.70	2 (16%)
14	CLA	M	156	-	23,32,73	7.53	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	M	157	-	23,32,73	7.52	5 (21%)	12,54,113	1.71	2 (16%)
14	CLA	M	158	-	23,32,73	7.58	5 (21%)	12,54,113	1.72	2 (16%)
14	CLA	M	159	-	23,32,73	7.57	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	M	160	-	23,32,73	7.60	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	M	161	-	23,32,73	7.62	5 (21%)	12,54,113	1.71	2 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
14	CLA	M	162	-	23,32,73	7.60	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	M	163	-	23,32,73	7.56	5 (21%)	12,54,113	1.68	2 (16%)
14	CLA	M	164	-	23,32,73	7.48	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	M	165	-	23,32,73	7.74	5 (21%)	12,54,113	1.68	2 (16%)
14	CLA	M	166	-	23,32,73	7.62	5 (21%)	12,54,113	1.68	2 (16%)
14	CLA	M	167	-	23,32,73	7.49	5 (21%)	12,54,113	1.69	2 (16%)
14	CLA	M	168	-	23,32,73	7.57	5 (21%)	12,54,113	1.70	2 (16%)
16	HEM	O	58	-	22,32,50	30.45	6 (27%)	16,54,82	1.51	3 (18%)
14	CLA	P	313	-	23,32,73	7.49	5 (21%)	12,54,113	1.70	2 (16%)
16	HEM	R	88	-	22,32,50	30.45	6 (27%)	16,54,82	1.53	3 (18%)

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
13	TYR	A	174	1	-	0/0/0/8	0/1/1/1
14	CLA	A	175	-	-	0/0/66/135	0/0/8/9
14	CLA	A	176	-	-	0/0/66/135	0/0/8/9
14	CLA	A	177	-	-	0/0/66/135	0/0/8/9
15	PHO	A	178	-	-	0/5/34/103	0/0/5/6
14	CLA	A	179	-	-	0/0/66/135	0/0/8/9
13	TYR	B	176	2	-	0/0/0/8	0/1/1/1
14	CLA	B	177	-	-	0/0/66/135	0/0/8/9
15	PHO	B	178	-	-	0/5/34/103	0/0/5/6
14	CLA	B	179	-	-	0/0/66/135	0/0/8/9
17	PLA	B	180	-	-	0/0/0/23	0/1/1/1
14	CLA	C	157	-	-	0/0/66/135	0/0/8/9
14	CLA	C	158	-	-	0/0/66/135	0/0/8/9
14	CLA	C	159	-	-	0/0/66/135	0/0/8/9
14	CLA	C	160	-	-	0/0/66/135	0/0/8/9
14	CLA	C	161	-	-	0/0/66/135	0/0/8/9
14	CLA	C	162	-	-	0/0/66/135	0/0/8/9
14	CLA	C	163	-	-	0/0/66/135	0/0/8/9
14	CLA	C	164	-	-	0/0/66/135	0/0/8/9
14	CLA	C	165	-	-	0/0/66/135	0/0/8/9
14	CLA	C	166	-	-	0/0/66/135	0/0/8/9
14	CLA	C	167	-	-	0/0/66/135	0/0/8/9
14	CLA	C	168	-	-	0/0/66/135	0/0/8/9

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
14	CLA	D	156	-	-	0/0/66/135	0/0/8/9
14	CLA	D	157	-	-	0/0/66/135	0/0/8/9
14	CLA	D	158	-	-	0/0/66/135	0/0/8/9
14	CLA	D	159	-	-	0/0/66/135	0/0/8/9
14	CLA	D	160	-	-	0/0/66/135	0/0/8/9
14	CLA	D	161	-	-	0/0/66/135	0/0/8/9
14	CLA	D	162	-	-	0/0/66/135	0/0/8/9
14	CLA	D	163	-	-	0/0/66/135	0/0/8/9
14	CLA	D	164	-	-	0/0/66/135	0/0/8/9
14	CLA	D	165	-	-	0/0/66/135	0/0/8/9
14	CLA	D	166	-	-	0/0/66/135	0/0/8/9
14	CLA	D	167	-	-	0/0/66/135	0/0/8/9
14	CLA	D	168	-	-	0/0/66/135	0/0/8/9
16	HEM	F	31	-	-	0/0/79/114	0/0/8/8
14	CLA	G	313	-	-	0/0/66/135	0/0/8/9
16	HEM	I	88	-	-	0/0/79/114	0/0/8/8
13	TYR	J	174	1	-	0/0/0/8	0/1/1/1
14	CLA	J	175	-	-	0/0/66/135	0/0/8/9
14	CLA	J	176	-	-	0/0/66/135	0/0/8/9
14	CLA	J	177	-	-	0/0/66/135	0/0/8/9
15	PHO	J	178	-	-	0/5/34/103	0/0/5/6
14	CLA	J	179	-	-	0/0/66/135	0/0/8/9
13	TYR	K	176	2	-	0/0/0/8	0/1/1/1
14	CLA	K	177	-	-	0/0/66/135	0/0/8/9
15	PHO	K	178	-	-	0/5/34/103	0/0/5/6
14	CLA	K	179	-	-	0/0/66/135	0/0/8/9
17	PLA	K	180	-	-	0/0/0/23	0/1/1/1
14	CLA	L	157	-	-	0/0/66/135	0/0/8/9
14	CLA	L	158	-	-	0/0/66/135	0/0/8/9
14	CLA	L	159	-	-	0/0/66/135	0/0/8/9
14	CLA	L	160	-	-	0/0/66/135	0/0/8/9
14	CLA	L	161	-	-	0/0/66/135	0/0/8/9
14	CLA	L	162	-	-	0/0/66/135	0/0/8/9
14	CLA	L	163	-	-	0/0/66/135	0/0/8/9
14	CLA	L	164	-	-	0/0/66/135	0/0/8/9
14	CLA	L	165	-	-	0/0/66/135	0/0/8/9
14	CLA	L	166	-	-	0/0/66/135	0/0/8/9
14	CLA	L	167	-	-	0/0/66/135	0/0/8/9
14	CLA	L	168	-	-	0/0/66/135	0/0/8/9
14	CLA	M	156	-	-	0/0/66/135	0/0/8/9
14	CLA	M	157	-	-	0/0/66/135	0/0/8/9
14	CLA	M	158	-	-	0/0/66/135	0/0/8/9

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
14	CLA	M	159	-	-	0/0/66/135	0/0/8/9
14	CLA	M	160	-	-	0/0/66/135	0/0/8/9
14	CLA	M	161	-	-	0/0/66/135	0/0/8/9
14	CLA	M	162	-	-	0/0/66/135	0/0/8/9
14	CLA	M	163	-	-	0/0/66/135	0/0/8/9
14	CLA	M	164	-	-	0/0/66/135	0/0/8/9
14	CLA	M	165	-	-	0/0/66/135	0/0/8/9
14	CLA	M	166	-	-	0/0/66/135	0/0/8/9
14	CLA	M	167	-	-	0/0/66/135	0/0/8/9
14	CLA	M	168	-	-	0/0/66/135	0/0/8/9
16	HEM	O	58	-	-	0/0/79/114	0/0/8/8
14	CLA	P	313	-	-	0/0/66/135	0/0/8/9
16	HEM	R	88	-	-	0/0/79/114	0/0/8/8

The worst 5 of 406 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	A	178	PHO	C3C-C4C	106.85	1.53	1.38
15	J	178	PHO	C3C-C4C	106.83	1.53	1.38
15	B	178	PHO	C3C-C4C	106.68	1.53	1.38
15	K	178	PHO	C3C-C4C	106.13	1.53	1.38
16	R	88	HEM	C2A-C1A	104.45	1.52	1.38

The worst 5 of 152 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	L	157	CLA	C4B-CHC-C1C	3.48	132.05	127.47
14	C	157	CLA	C4B-CHC-C1C	3.43	131.98	127.47
14	L	167	CLA	C4B-CHC-C1C	3.42	131.97	127.47
14	L	166	CLA	C4B-CHC-C1C	3.41	131.96	127.47
14	C	167	CLA	C4B-CHC-C1C	3.41	131.96	127.47

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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 will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.