



wwPDB X-ray Structure Validation Summary Report

Feb 28, 2014 – 06:05 PM GMT

PDB ID : 1JGX
Title : Photosynthetic Reaction Center Mutant With Thr M 21 Replaced With Asp
Authors : Camara-Artigas, A.; Magee, C.L.; Williams, J.C.; Allen, J.P.
Deposited on : 2001-06-27
Resolution : 3.01 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>

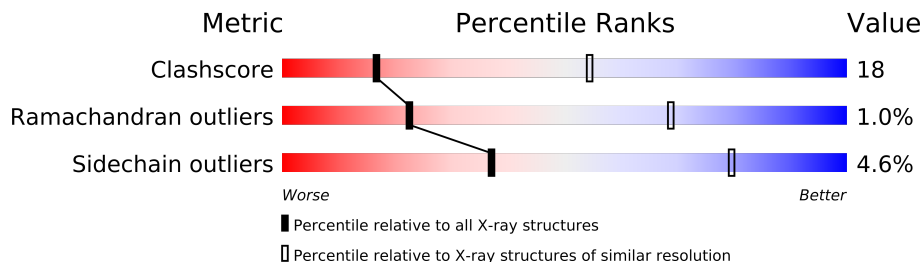
The following versions of software and data (see [references](#)) were used in the production of this report:

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 3.01 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	1732 (3.04-3.00)
Ramachandran outliers	78287	1669 (3.04-3.00)
Sidechain outliers	78261	1672 (3.04-3.00)

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 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	L	281	
2	M	307	
3	H	260	

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 7067 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 Photosynthetic Reaction Center L subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	281	Total	C	N	O	S	0	0	0
			2232	1507	355	362	8			

- Molecule 2 is a protein called Photosynthetic Reaction Center M subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	M	302	Total	C	N	O	S	0	0	0
			2409	1607	394	398	10			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
M	21	ASP	THR	ENGINEERED	UNP P02953

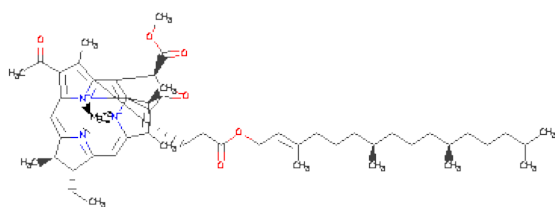
- Molecule 3 is a protein called Photosynthetic Reaction Center H subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	240	Total	C	N	O	S	0	0	0
			1829	1169	314	337	9			

- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe).

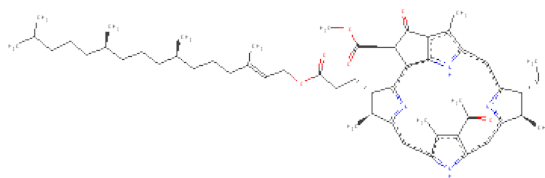
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	M	1	Total	Fe	0	0
			1	1		

- Molecule 5 is BACTERIOCHLOROPHYLL A (three-letter code: BCL) (formula: C₅₅H₇₄MgN₄O₆).



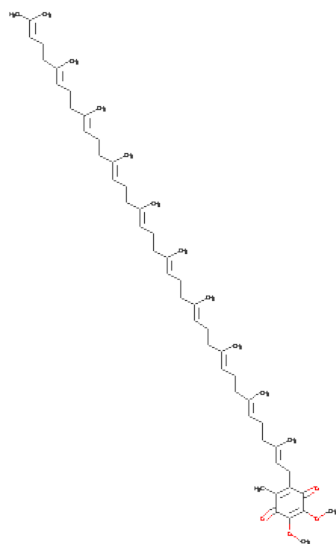
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
5	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
5	M	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
5	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

- Molecule 6 is BACTERIOPHEOPHYTIN A (three-letter code: BPH) (formula: $C_{55}H_{76}N_4O_6$).



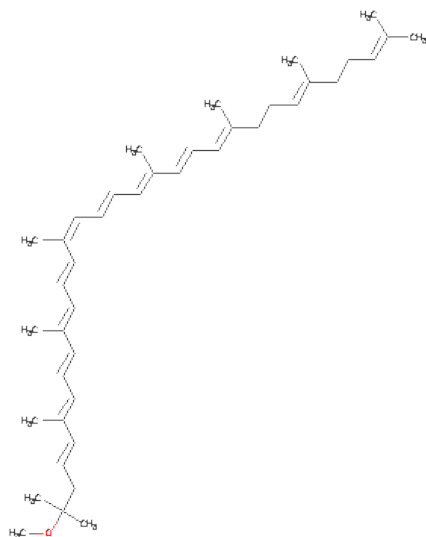
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	M	1	Total	C	N	O	0	0
			65	55	4	6		
6	L	1	Total	C	N	O	0	0
			65	55	4	6		

- Molecule 7 is UBIQUINONE-10 (three-letter code: U10) (formula: $C_{59}H_{90}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	M	1	Total	C	O	0	0
			48	44	4		

- Molecule 8 is SPHEROIDENE (three-letter code: SPO) (formula: $C_{41}H_{60}O$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	M	1	Total	C	O	0	0
			42	41	1		

- Molecule 9 is water.

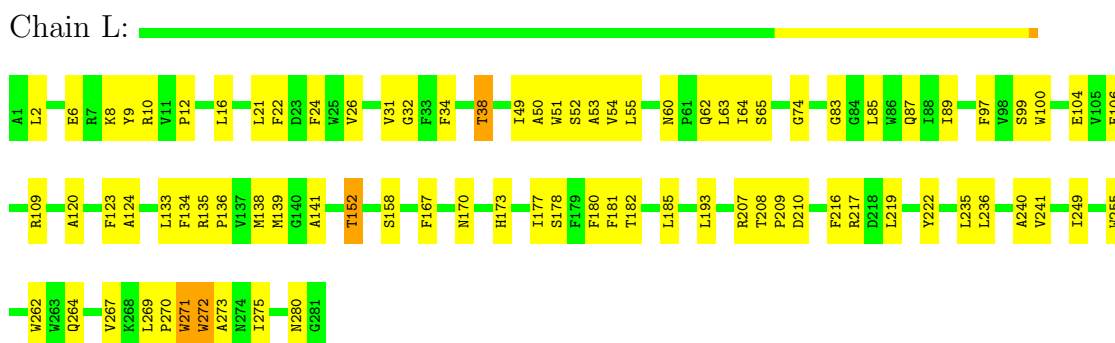
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	H	47	Total	O	0	0
			47	47		
9	L	28	Total	O	0	0
			28	28		
9	M	37	Total	O	0	0
			37	37		

3 Residue-property plots

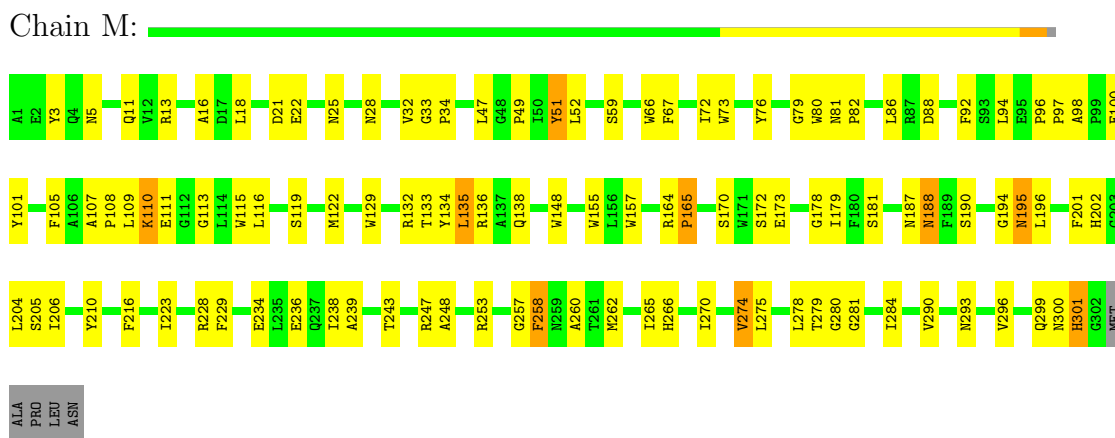
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

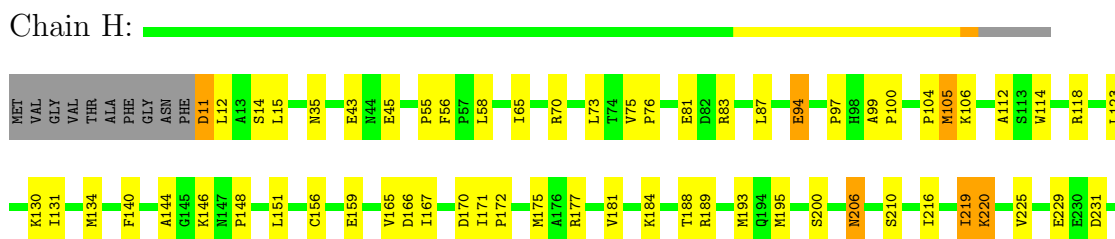
- Molecule 1: Photosynthetic Reaction Center L subunit



- Molecule 2: Photosynthetic Reaction Center M subunit



- Molecule 3: Photosynthetic Reaction Center H subunit



G235	
G239	
G240	
L241	
Y242	
Y243	
A244	
A245	
R248	
K249	
S250	
VAL	
VAL	
ALA	
ALA	
MET	
LEU	
ALA	
GLU	
TYR	
ALA	

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	142.11Å 142.11Å 187.74Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.97 – 3.01	Depositor
% Data completeness (in resolution range)	92.5 (29.97-3.01)	Depositor
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.211 , 0.249	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7067	wwPDB-VP
Average B, all atoms (Å ²)	50.0	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: BCL, BPH, U10, FE, SPO

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	L	0.41	0/2320	0.59	0/3175
2	M	0.42	0/2501	0.61	1/3414 (0.0%)
3	H	0.40	0/1877	0.67	0/2553
All	All	0.41	0/6698	0.62	1/9142 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	258	PHE	N-CA-CB	-5.04	101.53	110.60

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	L	2232	0	2187	79	0
2	M	2409	0	2318	110	0
3	H	1829	0	1836	57	0
4	M	1	0	0	0	0
5	L	198	0	222	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	M	66	0	74	7	0
6	L	65	0	76	7	0
6	M	65	0	76	9	0
7	M	48	0	63	5	0
8	M	42	0	60	1	0
9	H	47	0	0	2	0
9	L	28	0	0	4	0
9	M	37	0	0	18	0
All	All	7067	0	6912	248	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 18.

The worst 5 of 248 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:M:243:THR:CG2	2:M:247:ARG:HE	1.71	1.03
2:M:260:ALA:CB	9:M:1027:HOH:O	2.12	0.97
1:L:60:ASN:HD22	1:L:63:LEU:HG	1.27	0.96
2:M:260:ALA:HB3	9:M:1027:HOH:O	1.68	0.93
1:L:97:PHE:HA	9:L:1014:HOH:O	1.71	0.91

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	L	279/281 (99%)	257 (92%)	21 (8%)	1 (0%)	43	85
2	M	300/307 (98%)	270 (90%)	24 (8%)	6 (2%)	11	47
3	H	238/260 (92%)	220 (92%)	17 (7%)	1 (0%)	43	85
All	All	817/848 (96%)	747 (91%)	62 (8%)	8 (1%)	22	69

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	M	110	LYS
2	M	301	HIS
1	L	273	ALA
2	M	179	ILE
2	M	92	PHE

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	L	220/220 (100%)	209 (95%)	11 (5%)	34	76
2	M	236/240 (98%)	227 (96%)	9 (4%)	44	85
3	H	195/208 (94%)	185 (95%)	10 (5%)	33	75
All	All	651/668 (98%)	621 (95%)	30 (5%)	37	79

5 of 30 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	M	86	LEU
2	M	188	ASN
3	H	220	LYS
2	M	165	PRO
2	M	216	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
2	M	28	ASN
3	H	206	ASN
2	M	202	HIS
1	L	173	HIS
2	M	299	GLN

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 9 ligands modelled in this entry, 1 is 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
5	BCL	L	850	-	74,74,74	2.33	16 (21%)	97,115,115	5.14	30 (30%)
5	BCL	L	851	-	74,74,74	1.48	12 (16%)	97,115,115	1.92	23 (23%)
5	BCL	L	853	-	74,74,74	1.51	14 (18%)	97,115,115	2.18	27 (27%)
6	BPH	L	855	-	70,70,70	1.50	7 (10%)	94,101,101	2.35	23 (24%)
5	BCL	M	852	-	74,74,74	1.45	11 (14%)	97,115,115	1.95	25 (25%)
6	BPH	M	854	-	70,70,70	1.53	9 (12%)	94,101,101	1.95	22 (23%)
7	U10	M	857	-	48,48,63	2.18	15 (31%)	59,61,79	2.41	22 (37%)
8	SPO	M	859	-	41,41,41	3.32	25 (60%)	50,50,50	2.98	18 (36%)

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
5	BCL	L	850	-	-	0/41/137/137	0/0/9/9
5	BCL	L	851	-	-	0/41/137/137	0/0/9/9
5	BCL	L	853	-	-	0/41/137/137	0/0/9/9
6	BPH	L	855	-	2/2/18/22	0/49/105/105	0/0/6/6
5	BCL	M	852	-	-	0/41/137/137	0/0/9/9

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	BPH	M	854	-	2/2/18/22	0/49/105/105	0/0/6/6
7	U10	M	857	-	-	0/45/69/87	0/1/1/1
8	SPO	M	859	-	-	0/47/47/47	0/0/0/0

The worst 5 of 109 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	L	850	BCL	C3B-C4B	11.85	1.58	1.40
5	L	850	BCL	C1B-NB	-9.00	1.24	1.34
8	M	859	SPO	C15-C16	8.61	1.57	1.34
7	M	857	U10	C6-C1	6.95	1.52	1.35
6	M	854	BPH	C1D-CHD	6.80	1.43	1.35

The worst 5 of 190 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	L	850	BCL	OBB-CAB-C3B	28.54	162.90	120.07
5	L	850	BCL	C4B-C3B-C2B	-22.97	79.19	106.97
5	L	850	BCL	C3B-C4B-NB	16.98	125.15	108.64
5	L	850	BCL	OBB-CAB-CBB	-16.88	78.19	120.13
5	L	850	BCL	CBB-CAB-C3B	-12.18	84.71	120.30

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
6	L	855	BPH	C8
6	L	855	BPH	C13
6	M	854	BPH	C8
6	M	854	BPH	C13

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.