



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 01:14 am GMT

PDB ID : 3ZC1  
Title : Crystal structure of AfC3PO  
Authors : Parizotto, E.A.; Lowe, E.D.; Parker, J.S.  
Deposited on : 2012-11-14  
Resolution : 3.27 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

---

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

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

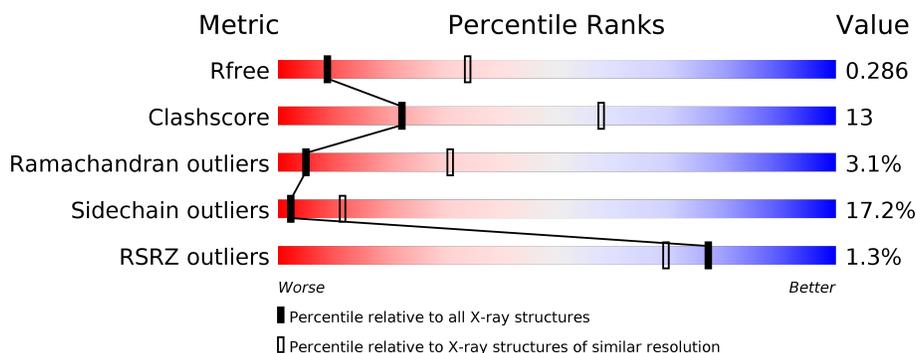
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.27 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))
$R_{free}$	100719	1852 (3.32-3.20)
Clashscore	112137	2036 (3.32-3.20)
Ramachandran outliers	110173	2000 (3.32-3.20)
Sidechain outliers	110143	1998 (3.32-3.20)
RSRZ outliers	101464	1861 (3.32-3.20)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	199	 0% (upper red bar), 64% (green), 24% (yellow), 7% (orange), 5% (grey)
1	B	199	 2% (upper red bar), 56% (green), 33% (yellow), 6% (orange), 5% (grey)
1	C	199	 0% (upper red bar), 58% (green), 30% (yellow), 7% (orange), 5% (grey)
1	D	199	 2% (upper red bar), 54% (green), 33% (yellow), 8% (orange), 6% (grey)
1	E	199	 0% (upper red bar), 55% (green), 31% (yellow), 10% (orange), 5% (grey)
1	F	199	 0% (upper red bar), 56% (green), 32% (yellow), 7% (orange), 5% (grey)

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
1	G	199	 <p>%</p> <p>56% 33% 6% • 5%</p>
1	H	199	 <p>3%</p> <p>59% 29% 6% 6%</p>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 11820 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 AFTRAX.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	190	1487	950	250	279	8	0	0	0
1	B	190	1495	950	255	282	8	0	0	0
1	C	189	1478	939	249	283	7	0	0	0
1	D	187	1479	943	249	280	7	0	0	0
1	E	190	1494	955	255	276	8	0	0	0
1	F	190	1486	948	251	279	8	0	0	0
1	G	189	1451	920	244	279	8	0	0	0
1	H	187	1437	911	245	273	8	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP O28024
A	-1	PRO	-	EXPRESSION TAG	UNP O28024
A	0	HIS	-	EXPRESSION TAG	UNP O28024
B	-2	GLY	-	EXPRESSION TAG	UNP O28024
B	-1	PRO	-	EXPRESSION TAG	UNP O28024
B	0	HIS	-	EXPRESSION TAG	UNP O28024
C	-2	GLY	-	EXPRESSION TAG	UNP O28024
C	-1	PRO	-	EXPRESSION TAG	UNP O28024
C	0	HIS	-	EXPRESSION TAG	UNP O28024
D	-2	GLY	-	EXPRESSION TAG	UNP O28024
D	-1	PRO	-	EXPRESSION TAG	UNP O28024
D	0	HIS	-	EXPRESSION TAG	UNP O28024
E	-2	GLY	-	EXPRESSION TAG	UNP O28024

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
E	-1	PRO	-	EXPRESSION TAG	UNP O28024
E	0	HIS	-	EXPRESSION TAG	UNP O28024
F	-2	GLY	-	EXPRESSION TAG	UNP O28024
F	-1	PRO	-	EXPRESSION TAG	UNP O28024
F	0	HIS	-	EXPRESSION TAG	UNP O28024
G	-2	GLY	-	EXPRESSION TAG	UNP O28024
G	-1	PRO	-	EXPRESSION TAG	UNP O28024
G	0	HIS	-	EXPRESSION TAG	UNP O28024
H	-2	GLY	-	EXPRESSION TAG	UNP O28024
H	-1	PRO	-	EXPRESSION TAG	UNP O28024
H	0	HIS	-	EXPRESSION TAG	UNP O28024

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

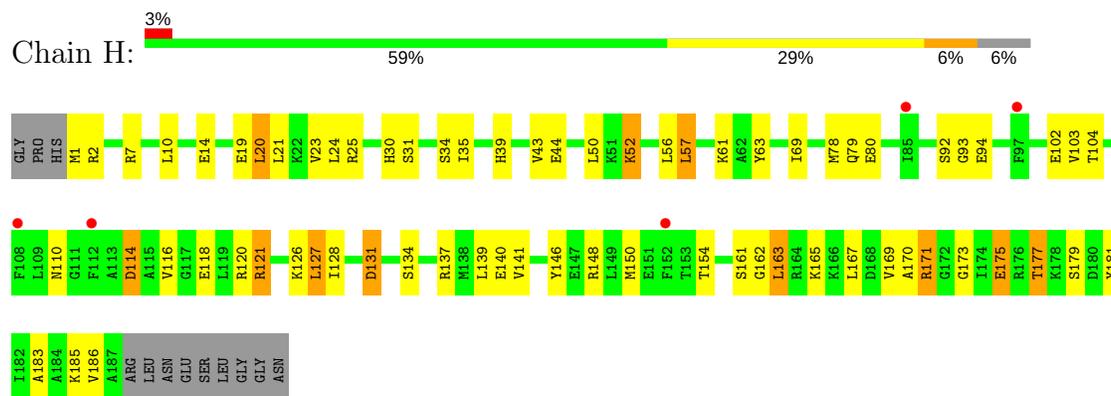
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	G	1	Total Mg 1 1	0	0
2	B	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0
2	F	1	Total Mg 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	3	Total O 3 3	0	0
3	D	1	Total O 1 1	0	0
3	F	3	Total O 3 3	0	0
3	G	1	Total O 1 1	0	0







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	183.31Å 183.31Å 111.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	53.56 – 3.27 70.75 – 3.27	Depositor EDS
% Data completeness (in resolution range)	99.9 (53.56-3.27) 98.1 (70.75-3.27)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.78 (at 3.26Å)	Xtrriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.221 , 0.281 0.224 , 0.286	Depositor DCC
$R_{free}$ test set	1984 reflections (6.72%)	DCC
Wilson B-factor (Å <sup>2</sup> )	102.8	Xtrriage
Anisotropy	0.318	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 72.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.25$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	11820	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	103.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.04% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 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: MG

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	0.49	0/1506	0.63	0/2025
1	B	0.48	0/1515	0.65	0/2038
1	C	0.49	0/1497	0.66	0/2016
1	D	0.47	0/1499	0.62	1/2016 (0.0%)
1	E	0.50	1/1513 (0.1%)	0.63	0/2031
1	F	0.50	0/1505	0.65	0/2024
1	G	0.49	0/1470	0.66	0/1983
1	H	0.48	0/1456	0.63	1/1963 (0.1%)
All	All	0.49	1/11961 (0.0%)	0.64	2/16096 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	6	CYS	CB-SG	5.27	1.91	1.82

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	D	57	LEU	CA-CB-CG	5.68	128.36	115.30
1	H	57	LEU	CA-CB-CG	5.56	128.09	115.30

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	1487	0	1485	33	0
1	B	1495	0	1479	45	0
1	C	1478	0	1454	47	0
1	D	1479	0	1475	48	0
1	E	1494	0	1513	48	0
1	F	1486	0	1486	46	0
1	G	1451	0	1395	48	0
1	H	1437	0	1387	33	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
3	B	3	0	0	0	0
3	D	1	0	0	0	0
3	F	3	0	0	0	0
3	G	1	0	0	0	0
All	All	11820	0	11674	309	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 13.

The worst 5 of 309 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)
1:G:2:ARG:NH1	1:G:6:CYS:SG	2.50	0.85
1:C:2:ARG:NH1	1:C:6:CYS:SG	2.54	0.80
1:F:161:SER:OG	1:F:162:GLY:N	2.20	0.74
1:D:183:ALA:O	1:D:185:LYS:N	2.22	0.72
1:H:183:ALA:O	1:H:185:LYS:N	2.22	0.72

There are no symmetry-related clashes.

## 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	Percentiles	
1	A	188/199 (94%)	164 (87%)	15 (8%)	9 (5%)	2	19
1	B	188/199 (94%)	159 (85%)	25 (13%)	4 (2%)	8	41
1	C	187/199 (94%)	160 (86%)	23 (12%)	4 (2%)	8	41
1	D	185/199 (93%)	152 (82%)	29 (16%)	4 (2%)	8	39
1	E	188/199 (94%)	164 (87%)	16 (8%)	8 (4%)	3	22
1	F	188/199 (94%)	164 (87%)	20 (11%)	4 (2%)	8	41
1	G	187/199 (94%)	159 (85%)	23 (12%)	5 (3%)	6	34
1	H	185/199 (93%)	147 (80%)	30 (16%)	8 (4%)	3	22
All	All	1496/1592 (94%)	1269 (85%)	181 (12%)	46 (3%)	5	31

5 of 46 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	ARG
1	B	2	ARG
1	B	42	LYS
1	B	131	ASP
1	E	2	ARG

### 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	Percentiles	
1	A	148/167 (89%)	123 (83%)	25 (17%)	2	12
1	B	149/167 (89%)	124 (83%)	25 (17%)	2	12
1	C	147/167 (88%)	127 (86%)	20 (14%)	4	20
1	D	149/167 (89%)	120 (80%)	29 (20%)	1	8
1	E	150/167 (90%)	125 (83%)	25 (17%)	2	12
1	F	149/167 (89%)	120 (80%)	29 (20%)	1	8

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	G	140/167 (84%)	119 (85%)	21 (15%)	3 16
1	H	138/167 (83%)	111 (80%)	27 (20%)	1 8
All	All	1170/1336 (88%)	969 (83%)	201 (17%)	2 11

5 of 201 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	141	VAL
1	E	99	ILE
1	H	56	LEU
1	D	164	ARG
1	E	20	LEU

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

Mol	Chain	Res	Type
1	D	79	GLN
1	H	79	GLN
1	E	39	HIS
1	B	79	GLN
1	F	110	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](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	190/199 (95%)	-0.01	2 (1%) 80 73	70, 100, 132, 157	0
1	B	190/199 (95%)	-0.08	4 (2%) 64 54	61, 99, 141, 157	0
1	C	189/199 (94%)	-0.09	1 (0%) 90 87	68, 97, 134, 159	0
1	D	187/199 (93%)	-0.05	3 (1%) 72 63	75, 104, 137, 157	0
1	E	190/199 (95%)	-0.10	1 (0%) 90 87	69, 100, 131, 157	0
1	F	190/199 (95%)	-0.08	2 (1%) 80 73	57, 97, 138, 151	0
1	G	189/199 (94%)	-0.08	1 (0%) 90 87	70, 99, 139, 160	0
1	H	187/199 (93%)	0.09	5 (2%) 55 46	75, 102, 137, 157	0
All	All	1512/1592 (94%)	-0.05	19 (1%) 77 69	57, 100, 139, 160	0

The worst 5 of 19 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	97	PHE	4.5
1	B	69	ILE	4.4
1	H	152	PHE	3.5
1	H	85	ILE	3.4
1	H	112	PHE	3.2

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

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	MG	B	197	1/1	0.92	0.07	-	93,93,93,93	0
2	MG	C	197	1/1	0.90	0.18	-	75,75,75,75	0
2	MG	F	197	1/1	0.91	0.12	-	95,95,95,95	0
2	MG	D	197	1/1	0.94	0.07	-	90,90,90,90	0
2	MG	G	197	1/1	0.94	0.10	-	74,74,74,74	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.