



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 1RJO
Title : AGAO + Xe
Authors : Guss, J.M.; Trambaiolo, D.M.; Duff, A.P.
Deposited on : 2003-11-19
Resolution : 1.67 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.1

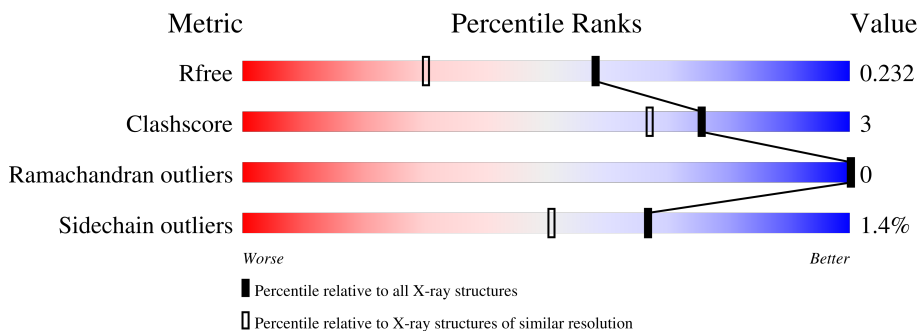
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 1.67 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}	130704	6780 (1.70-1.66)
Clashscore	141614	7310 (1.70-1.66)
Ramachandran outliers	138981	7173 (1.70-1.66)
Sidechain outliers	138945	7172 (1.70-1.66)

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	646	 91% 5%

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
5	XE	A	901	-	-	X	-
5	XE	A	902	-	-	X	-
5	XE	A	903	-	-	X	-
5	XE	A	904	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	XE	A	905	-	-	X	-
5	XE	A	909	-	-	X	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 5463 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 Phenylethylamine oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	620	4884	3087	855	932	10	0	4	0

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	382	TPQ	TYR	modified residue	UNP P46881
A	639	SER	-	cloning artifact	UNP P46881
A	640	ASN	-	cloning artifact	UNP P46881
A	641	TRP	-	cloning artifact	UNP P46881
A	642	SER	-	cloning artifact	UNP P46881
A	643	HIS	-	cloning artifact	UNP P46881
A	644	PRO	-	cloning artifact	UNP P46881
A	645	GLN	-	cloning artifact	UNP P46881
A	646	PHE	-	cloning artifact	UNP P46881
A	647	GLU	-	cloning artifact	UNP P46881
A	648	LYS	-	cloning artifact	UNP P46881

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Cu	0	0
			1	1		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0

- Molecule 5 is XENON (three-letter code: XE) (formula: Xe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	7	Total Xe 7 7	0	0

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 6 3 3	0	0
6	A	1	Total C O 6 3 3	0	0
6	A	1	Total C O 10 6 4	0	1

- Molecule 7 is water.

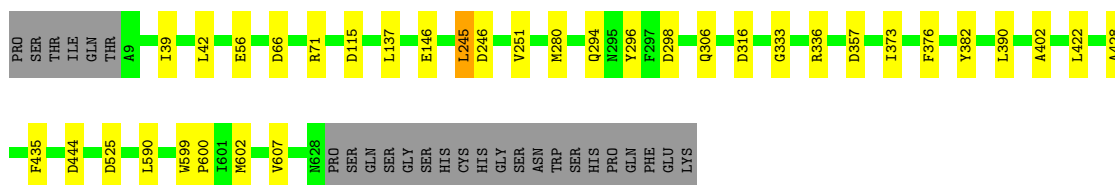
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	543	Total O 543 543	0	0

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.

- Molecule 1: Phenylethylamine oxidase

Chain A:  91% 5%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	158.01Å 63.19Å 92.15Å 90.00° 112.09° 90.00°	Depositor
Resolution (Å)	26.30 – 1.67 85.38 – 1.67	Depositor EDS
% Data completeness (in resolution range)	97.2 (26.30-1.67) 97.1 (85.38-1.67)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 1.67Å)	Xtrriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.157 , 0.178 0.217 , 0.232	Depositor DCC
R_{free} test set	7723 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	22.5	Xtrriage
Anisotropy	0.225	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 54.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5463	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: XE, GOL, SO4, TPQ, NA, CU

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	0.36	0/5008	0.73	7/6818 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	298	ASP	CB-CG-OD2	5.60	123.34	118.30
1	A	66	ASP	CB-CG-OD2	5.55	123.29	118.30
1	A	115	ASP	CB-CG-OD2	5.45	123.20	118.30
1	A	357	ASP	CB-CG-OD2	5.42	123.17	118.30
1	A	316	ASP	CB-CG-OD2	5.34	123.11	118.30
1	A	246	ASP	CB-CG-OD2	5.22	123.00	118.30
1	A	525	ASP	CB-CG-OD2	5.17	122.95	118.30

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	4884	0	4705	30	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	7	0	0	21	0
6	A	22	0	28	0	0
7	A	543	0	0	0	0
All	All	5463	0	4733	30	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 3.

All (30) 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:A:590[A]:LEU:HD12	5:A:902:XE:XE	1.93	1.45
1:A:590[A]:LEU:CD1	5:A:902:XE:XE	2.49	1.38
1:A:590[A]:LEU:HD11	5:A:902:XE:XE	2.24	1.14
1:A:245:LEU:HD22	5:A:901:XE:XE	2.45	0.94
1:A:245:LEU:CD2	5:A:901:XE:XE	2.97	0.90
1:A:390:LEU:HG	5:A:904:XE:XE	2.64	0.76
1:A:373:ILE:HD12	5:A:904:XE:XE	2.66	0.74
1:A:402:ALA:HB2	5:A:902:XE:XE	2.67	0.73
1:A:42:LEU:CD1	5:A:905:XE:XE	3.19	0.69
1:A:245:LEU:HD21	5:A:901:XE:XE	2.80	0.59
1:A:599:TRP:CD2	1:A:600:PRO:HA	2.38	0.59
1:A:590[B]:LEU:HD13	1:A:607:VAL:HB	1.92	0.52
1:A:280[A]:MET:HG3	1:A:435:PHE:CD2	2.45	0.52
1:A:590[B]:LEU:HD22	5:A:902:XE:XE	2.88	0.51
1:A:42:LEU:HD13	5:A:905:XE:XE	2.88	0.50
1:A:422:LEU:HD11	1:A:428:ALA:HB2	1.94	0.50
1:A:251:VAL:HG22	1:A:306:GLN:HB3	1.93	0.49
1:A:402:ALA:CB	5:A:902:XE:XE	3.39	0.49
1:A:71:ARG:HD3	5:A:903:XE:XE	2.91	0.48
1:A:294:GLN:HG3	1:A:296:TYR:CZ	2.48	0.48
1:A:280[A]:MET:HG3	1:A:435:PHE:CE2	2.49	0.47
1:A:607:VAL:HG12	5:A:902:XE:XE	2.94	0.46
1:A:71:ARG:CD	5:A:903:XE:XE	3.41	0.46
1:A:373:ILE:CD1	5:A:904:XE:XE	3.38	0.46
1:A:137:LEU:HD22	1:A:296:TYR:CZ	2.52	0.45
1:A:42:LEU:HD23	5:A:909:XE:XE	2.96	0.44
1:A:39:ILE:HD13	5:A:905:XE:XE	2.96	0.44
1:A:333:GLY:HA2	5:A:905:XE:XE	2.97	0.43
1:A:42:LEU:CD2	5:A:909:XE:XE	3.47	0.40

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	621/646 (96%)	602 (97%)	19 (3%)	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	Percentiles
1	A	517/537 (96%)	510 (99%)	7 (1%)	67 51

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	GLU
1	A	146	GLU
1	A	245	LEU
1	A	336	ARG
1	A	376	PHE
1	A	444	ASP
1	A	602	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	507	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](#)

1 non-standard protein/DNA/RNA residue is 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$
1	TPQ	A	382	1	13,14,15	1.87	4 (30%)	15,19,21	1.23	2 (13%)

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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	TPQ	A	382	1	-	1/5/22/24	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	382	TPQ	O5-C5	3.56	1.34	1.24
1	A	382	TPQ	O2-C2	3.36	1.33	1.24
1	A	382	TPQ	C3-C4	2.92	1.40	1.35
1	A	382	TPQ	C6-C1	2.05	1.40	1.34

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	382	TPQ	CB-CA-C	-3.59	104.75	111.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	382	TPQ	C6-C1-C2	2.03	120.20	118.64

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	382	TPQ	O-C-CA-CB

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 14 ligands modelled in this entry, 9 are monoatomic - leaving 5 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	$\# Z > 2$
6	GOL	A	1203[A]	-	5,5,5	0.41	0	5,5,5	0.21	0
4	SO4	A	1101	-	4,4,4	0.12	0	6,6,6	0.18	0
6	GOL	A	1201	-	5,5,5	0.32	0	5,5,5	0.19	0
6	GOL	A	1202	-	5,5,5	0.37	0	5,5,5	0.19	0
6	GOL	A	1203[B]	-	5,5,5	0.37	0	5,5,5	0.13	0

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
6	GOL	A	1203[A]	-	-	4/4/4/4	-
6	GOL	A	1202	-	-	0/4/4/4	-
6	GOL	A	1201	-	-	0/4/4/4	-
6	GOL	A	1203[B]	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1203[A]	GOL	O1-C1-C2-O2
6	A	1203[A]	GOL	O1-C1-C2-C3
6	A	1203[A]	GOL	O2-C2-C3-O3
6	A	1203[B]	GOL	O1-C1-C2-C3
6	A	1203[A]	GOL	C1-C2-C3-O3
6	A	1203[B]	GOL	O1-C1-C2-O2

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

6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.