

Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2022 – 06:18 pm GMT

PDB ID 6RGM

Title urate oxidase under 130 bar of krypton Authors Prange, T.; Colloc'h, N.; Carpentier, P.

2019-04-17 Deposited on

1.50 Å(reported) Resolution

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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity 4.02b-467

> 1.8.4, CSD as541be (2020) Mogul

Xtriage (Phenix) 1.13

EDS 2.26

Percentile statistics 20191225.v01 (using entries in the PDB archive December 25th 2019)

> Refmac 5.8.0267

CCP4 7.1.010 (Gargrove)

Ideal geometry (proteins) Engh & Huber (2001) Ideal geometry (DNA, RNA) Parkinson et al. (1996)

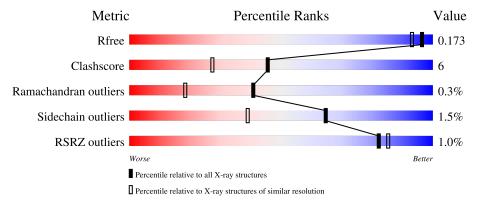
Validation Pipeline (wwPDB-VP) 2.26

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\ DIFFRACTION$

The reported resolution of this entry is 1.50 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.



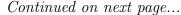
Metric	$\begin{array}{c} \textbf{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\AA)}) \end{array}$
R_{free}	130704	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (1.50-1.50)

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 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 <=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	302	91%	6% ••					

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
2	MPD	A	1103	-	-	X	-





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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ACT	A	1106	-	-	X	-
5	KR	A	1109	-	-	X	-



2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 3038 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.

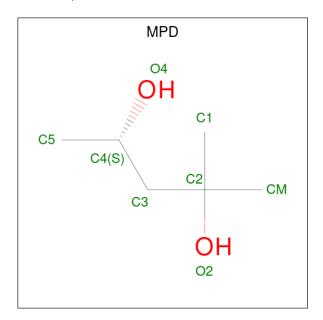
• Molecule 1 is a protein called Uricase.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	A	297	Total 2456	C 1554	N 424	O 469	S 9	0	16	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1000	ACE	-	expression tag	UNP Q00511

• Molecule 2 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



I	Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
	2	A	1	Total C O 8 6 2	0	0
	2	A	1	Total C O 8 6 2	0	0

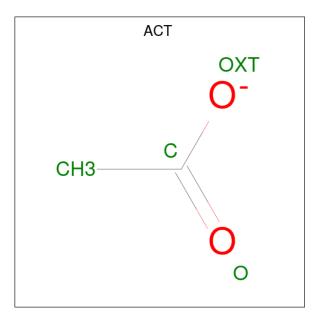
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total C (8 6 5) 2	0	0

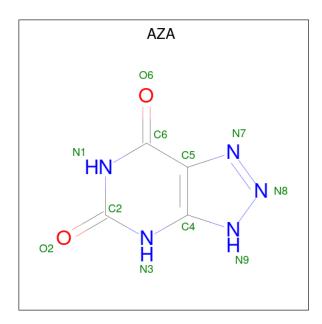
 \bullet Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $\mathrm{C_2H_3O_2}).$



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0

 \bullet Molecule 4 is 8-AZAXANTHINE (three-letter code: AZA) (formula: $\mathrm{C_4H_3N_5O_2}).$





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total 11	C 4	N 5	O 2	0	0

• Molecule 5 is KRYPTON (three-letter code: KR) (formula: Kr).

Mo	ol	Chain	Residues	Ator	ns	ZeroOcc	AltConf
5		A	3	Total 3	Kr 3	0	0

• Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

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

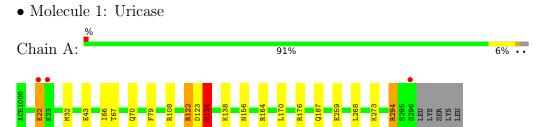
• Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	511	Total O 531 531	0	21



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 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.





4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants	78.25Å 94.94Å 104.08Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	70.14 - 1.50	Depositor
Resolution (A)	70.14 - 1.50	EDS
% Data completeness	99.9 (70.14-1.50)	Depositor
(in resolution range)	99.9 (70.14-1.50)	EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	3.71 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
P. P.	0.140 , 0.169	Depositor
R, R_{free}	0.145 , 0.173	DCC
R_{free} test set	3146 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	20.1	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	(Not available), (Not available)	EDS
L-test for twinning ²	$< L > = 0.49, < L^2> = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3038	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.54% of the height of the origin peak. No significant pseudotranslation is detected.

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



¹Intensities estimated from amplitudes.

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: NA, AZA, KR, ACE, ACT, MPD

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).

Mal	Chain	Bond lengths		Bond angles	
Mol		RMSZ	# Z > 5	RMSZ	# Z > 5
1	A	0.76	3/2548 (0.1%)	0.95	7/3452 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a maintain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(A)
1	A	259	GLU	CD-OE1	-5.57	1.19	1.25
1	A	22[A]	GLU	C-O	5.14	1.33	1.23
1	A	22[B]	GLU	C-O	5.14	1.33	1.23

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	123[A]	ASP	C-N-CA	6.40	137.69	121.70
1	A	123[B]	ASP	C-N-CA	6.40	137.69	121.70
1	A	108	ARG	NE-CZ-NH2	-5.58	117.51	120.30
1	A	22[A]	GLU	CA-C-N	-5.36	105.41	117.20
1	A	22[B]	GLU	CA-C-N	-5.36	105.41	117.20
1	A	122	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	A	294	ARG	NE-CZ-NH1	-5.03	117.78	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	A	124[A]	SER	Mainchain

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	2456	0	2446	20	0
2	A	24	0	42	15	0
3	A	12	0	9	5	0
4	A	11	0	3	1	0
5	A	3	0	0	2	0
6	A	1	0	0	0	0
7	A	531	0	0	7	0
All	All	3038	0	2500	32	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 6.

All (32) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
1100111 1	1100111 2	${ m distance} ({ m \AA})$	overlap (Å)
1:A:124[A]:SER:HB2	7:A:1505[A]:HOH:O	1.37	1.24
1:A:43:GLU:HG3	3:A:1106:ACT:H2	1.23	1.15
2:A:1102:MPD:HM1	2:A:1102:MPD:H52	1.48	0.94
1:A:66:ILE:HG22	2:A:1103:MPD:H11	1.47	0.93
2:A:1103:MPD:HM1	2:A:1103:MPD:H52	1.58	0.84
2:A:1102:MPD:H53	7:A:1247:HOH:O	1.85	0.76
1:A:176:ARG:HH12	4:A:1107:AZA:HN3	1.31	0.76
2:A:1102:MPD:H52	2:A:1102:MPD:CM	2.18	0.74
1:A:43:GLU:HG3	3:A:1106:ACT:CH3	2.13	0.73
2:A:1103:MPD:HM1	2:A:1103:MPD:C5	2.20	0.71
3:A:1106:ACT:OXT	7:A:1201:HOH:O	2.11	0.68
1:A:67:THR:OG1	2:A:1103:MPD:H13	1.94	0.66
1:A:32[A]:MET:HG2	5:A:1109:KR:KR	2.58	0.63
1:A:32[A]:MET:HE3	1:A:79:PHE:CE2	2.35	0.61
1:A:43:GLU:CG	3:A:1106:ACT:H2	2.15	0.61
1:A:164[B]:ARG:NH1	1:A:170:LEU:O	2.31	0.61

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Atom-1	Atom-2	Interatomic	Clash
7100111-1	1100111-2	${ m distance}({ m \AA})$	overlap (Å)
3:A:1106:ACT:C	7:A:1215:HOH:O	2.49	0.60
1:A:66:ILE:CG2	2:A:1103:MPD:H11	2.29	0.54
1:A:70:GLN:OE1	2:A:1103:MPD:H31	2.11	0.51
2:A:1102:MPD:H13	7:A:1632:HOH:O	2.12	0.48
2:A:1101:MPD:H51	7:A:1638:HOH:O	2.17	0.45
1:A:294:ARG:HH11	1:A:294:ARG:HG3	1.82	0.45
2:A:1101:MPD:HM3	2:A:1101:MPD:O4	2.18	0.44
1:A:32[A]:MET:CE	5:A:1109:KR:KR	3.27	0.43
1:A:294:ARG:HG3	1:A:294:ARG:NH1	2.34	0.43
1:A:67:THR:CA	2:A:1103:MPD:H13	2.49	0.42
1:A:32[B]:MET:CE	7:A:1277:HOH:O	2.67	0.42
2:A:1102:MPD:CM	2:A:1102:MPD:C5	2.91	0.41
1:A:22[A]:GLU:HA	1:A:22[A]:GLU:OE2	2.21	0.41
1:A:187[B]:GLN:HE21	1:A:187[B]:GLN:HB3	1.78	0.40
2:A:1101:MPD:O4	2:A:1101:MPD:CM	2.70	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	311/302 (103%)	304 (98%)	5 (2%)	2 (1%)	25 7

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124[A]	SER
1	A	124[B]	SER



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	278/267 (104%)	273 (98%)	5 (2%)	59 30	

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	122	ARG
1	A	156	ASN
1	A	268	LEU
1	A	273[A]	LYS
1	A	273[B]	LYS

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	156	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	Tuno	Chain	Res	Link	В	ond len	gths	Bond angles		
IVIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MPD	A	1102	-	7,7,7	0.27	0	9,10,10	0.53	0
2	MPD	A	1103	-	7,7,7	0.32	0	9,10,10	0.86	0
3	ACT	A	1105	-	1,3,3	3.59	1 (100%)	0,3,3	_	-
4	AZA	A	1107	-	9,12,12	0.95	1 (11%)	4,17,17	9.03	2 (50%)
3	ACT	A	1106	-	1,3,3	0.42	0	0,3,3	-	-
3	ACT	A	1104	-	1,3,3	4.37	1 (100%)	0,3,3	_	-
2	MPD	A	1101	-	7,7,7	0.19	0	9,10,10	0.51	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
2	MPD	A	1102	-	-	1/5/5/5	-
2	MPD	A	1103	-	=	1/5/5/5	-
4	AZA	A	1107	-	-	-	0/2/2/2
2	MPD	A	1101	-	=	0/5/5/5	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(\text{\AA})$
3	A	1104	ACT	СН3-С	4.37	1.54	1.48
3	A	1105	ACT	СН3-С	3.59	1.53	1.48
4	A	1107	AZA	C6-N1	2.48	1.37	1.33

All (2) bond angle outliers are listed below:

	Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\mathbf{Ideal}(^{o})$
Ī	4	A	1107	AZA	C2-N1-C6	15.88	128.55	115.14
Ī	4	A	1107	AZA	C5-C6-N1	-8.34	112.02	123.43

There are no chirality outliers.



All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1102	MPD	C2-C3-C4-O4
2	A	1103	MPD	C2-C3-C4-C5

There are no ring outliers.

5 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1102	MPD	5	0
2	A	1103	MPD	7	0
4	A	1107	AZA	1	0
3	A	1106	ACT	5	0
2	A	1101	MPD	3	0

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^{th} 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>	#RS	$\mathrm{SRZ}{>}2$	$OWAB(Å^2)$	Q<0.9
1	A	296/302 (98%)	-0.40	3 (1%)	82 85	14, 20, 36, 71	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	296	SER	4.8
1	A	22[A]	GLU	2.6
1	A	23[A]	LYS	2.5

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 monosaccharides 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. The B-factors column lists the minimum, median, 95^{th} 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	${f B-factors}({f \AA}^2)$	Q<0.9
2	MPD	A	1103	8/8	0.67	0.23	38,43,46,53	8
3	ACT	A	1104	4/4	0.72	0.18	35,42,45,48	0
2	MPD	A	1102	8/8	0.73	0.20	33,39,45,51	8
3	ACT	A	1105	4/4	0.79	0.14	29,39,42,53	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	${f B-factors}({f \AA}^2)$	Q<0.9
2	MPD	A	1101	8/8	0.81	0.18	43,46,58,61	8
3	ACT	A	1106	4/4	0.87	0.34	30,30,30,30	0
5	KR	A	1110	1/1	0.93	0.34	32,32,32,32	1
4	AZA	A	1107	11/11	0.97	0.06	15,15,17,18	0
6	NA	A	1111	1/1	0.98	0.06	21,21,21,21	0
5	KR	A	1108	1/1	0.99	0.12	28,28,28,28	0
5	KR	A	1109	1/1	1.00	0.25	22,22,22,22	1

6.5 Other polymers (i)

There are no such residues in this entry.

