

# Full wwPDB X-ray Structure Validation Report (i)

#### Mar 9, 2024 – 02:31 PM EST

PDB ID : 3S57

Title: ABH2 cross-linked with undamaged dsDNA-1 containing cofactors

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Deposited on : 2011-05-20

Resolution : 1.60 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 (i) 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 (1)) 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

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)

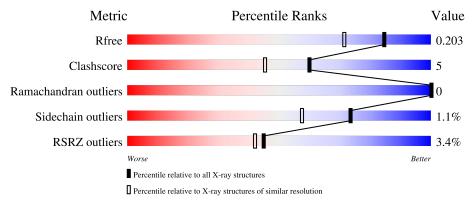
Validation Pipeline (wwPDB-VP) : 2.36

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\#  ext{Entries},  ext{ resolution range}( ext{Å}))$
$R_{free}$	130704	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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	204	89%	10% •				
2	В	14	79%	21%				
3	С	14	57%	36% 7%				



# 2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 2658 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 Alpha-ketoglutarate-dependent dioxygenase alkB homolog 2.

$\mathbf{Mol}$	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	A	204	Total 1727	C 1097	N 316	O 310	S 4	9	10	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference	
A	55	MET	-	expression tag	UNP Q6NS38	
A	67	SER	CYS	engineered mutation	UNP Q6NS38	
A	165	SER	CYS	engineered mutation	UNP Q6NS38	
A	169	CYS	GLY	engineered mutation	UNP Q6NS38	
A	192	SER	CYS	engineered mutation	UNP Q6NS38	

• Molecule 2 is a DNA chain called 5'-D(\*CP\*TP\*GP\*TP\*CP\*AP\*TP\*CP\*AP\*CP\*TP\*GP \*CP\*G)-3'.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	В	14	Total 280	C 135	N 48	O 84	P 13	0	0	0

• Molecule 3 is a DNA chain called 5'-D(\*TP\*CP\*GP\*CP\*AP\*GP\*TP\*GP\*AP\*TP\*GP\*A P\*CP\*A)-3'.

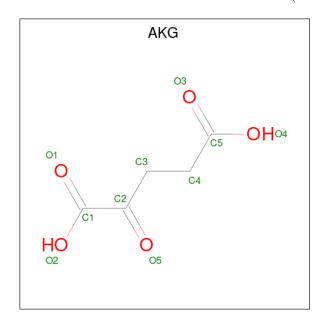
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
3	С	14	Total 286	C 137	N 55	O 81	P 13	0	0	0

• Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Mn 1 1	0	0

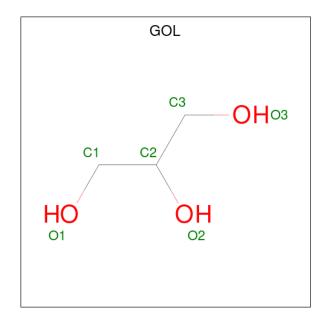


 $\bullet$  Molecule 5 is 2-OXOGLUTARIC ACID (three-letter code: AKG) (formula:  $C_5H_6O_5$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total 10	C 5	O 5	0	0

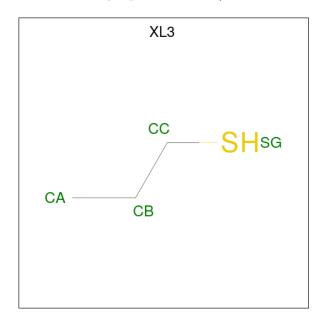
• Molecule 6 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



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



• Molecule 7 is propane-1-thiol (three-letter code: XL3) (formula:  $C_3H_8S$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	В	1	Total 4	C 3	S 1	0	0

• Molecule 8 is water.

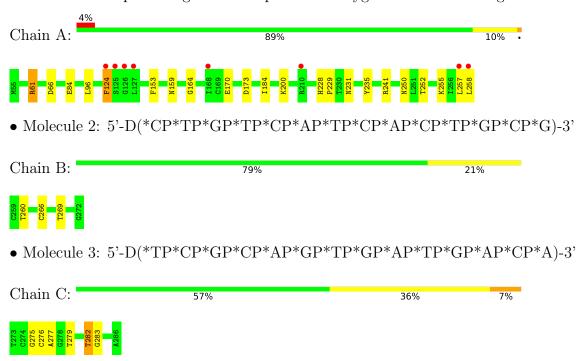
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	249	Total O 249 249	0	0
8	В	47	Total O 47 47	0	0
8	С	42	Total O 42 42	0	0



# 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 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.

• Molecule 1: Alpha-ketoglutarate-dependent dioxygenase alkB homolog 2





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	45.97Å 60.54Å 65.45Å	Denogitor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.85^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	20.00 - 1.60	Depositor
rtesolution (A)	32.11 - 1.49	EDS
% Data completeness	97.0 (20.00-1.60)	Depositor
(in resolution range)	96.6 (32.11-1.49)	EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	4.55  (at  1.50Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
P. P.	0.175 , $0.202$	Depositor
$R, R_{free}$	0.174 , $0.203$	DCC
$R_{free}$ test set	2876 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	19.2	Xtriage
Anisotropy	0.254	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.39 , 47.9	EDS
L-test for twinning <sup>2</sup>	$< L > = 0.50, < L^2> = 0.33$	Xtriage
Estimated twinning fraction	0.025 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	2658	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.60% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical 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.



<sup>&</sup>lt;sup>1</sup>Intensities 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: GOL, XL3, MN, AKG

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	Bo	nd lengths	Bond angles		
Mol Chain		RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.52	2/1774~(0.1%)	0.65	0/2400	
2	В	0.80	0/312	1.52	6/479 (1.3%)	
3	С	0.80	0/321	1.65	7/494 (1.4%)	
All	All	0.60	$2/2407 \; (0.1\%)$	1.01	$13/3373 \ (0.4\%)$	

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	Observed(A)	$Ideal(\AA)$
1	A	84[A]	GLU	CD-OE2	-5.43	1.19	1.25
1	A	84[B]	GLU	CD-OE2	-5.43	1.19	1.25

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$Ideal(^{o})$
3	С	279	DT	O4'-C4'-C3'	-10.28	99.83	106.00
3	С	275	DG	O4'-C1'-N9	9.15	114.41	108.00
3	С	277	DA	O4'-C1'-N9	-8.07	102.35	108.00
2	В	266	DC	O4'-C1'-N1	-6.89	103.18	108.00
2	В	260	DT	O4'-C1'-N1	-6.52	103.44	108.00
2	В	266	DC	P-O3'-C3'	6.35	127.32	119.70
3	С	282	DT	C1'-O4'-C4'	-6.22	103.88	110.10
3	С	276	DC	O4'-C1'-N1	-6.18	103.68	108.00
2	В	269	DT	C4-C5-C7	6.03	122.62	119.00
3	С	279	DT	C1'-O4'-C4'	-5.81	104.29	110.10
2	В	269	DT	C6-C5-C7	-5.62	119.53	122.90
3	С	279	DT	C4'-C3'-C2'	-5.03	98.57	103.10
2	В	266	DC	C5'-C4'-C3'	5.00	123.11	114.10

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	1727	0	1693	19	0
2	В	280	0	159	0	0
3	С	286	0	159	1	0
4	A	1	0	0	0	0
5	A	10	0	4	0	0
6	A	12	0	16	0	0
7	В	4	0	6	0	0
8	A	249	0	0	6	1
8	В	47	0	0	0	0
8	C	42	0	0	0	0
All	All	2658	0	2037	20	1

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 5.

All (20) 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:252[B]:THR:HG21	8:A:345:HOH:O	1.76	0.84
1:A:173[B]:ASP:H	1:A:231:ASN:HD21	1.29	0.79
1:A:173[A]:ASP:H	1:A:231:ASN:HD21	1.30	0.77
1:A:250[B]:ASN:ND2	8:A:16:HOH:O	1.71	0.68
1:A:252[B]:THR:CG2	8:A:345:HOH:O	2.39	0.65
1:A:96:LEU:HD22	1:A:124[A]:PHE:HD2	1.60	0.65
1:A:159:ASN:HB2	1:A:250[B]:ASN:HD22	1.69	0.56
1:A:257:LEU:O	1:A:258:LEU:HB2	2.07	0.54
1:A:96:LEU:HD22	1:A:124[A]:PHE:CD2	2.44	0.50
1:A:228[B]:HIS:CD2	1:A:229:PRO:HA	2.47	0.49
1:A:61:ARG:NH2	1:A:66:ASP:HB2	2.28	0.48
1:A:164:GLY:HA3	1:A:241:ARG:O	2.14	0.47
1:A:170[A]:GLU:HG3	1:A:235:TYR:HB3	1.97	0.47
1:A:170[B]:GLU:HG2	8:A:30:HOH:O	2.15	0.45
1:A:153:PHE:HA	1:A:255[A]:LYS:HD3	2.00	0.44
3:C:282:DT:H2"	3:C:283:DG:C8	2.53	0.44
1:A:170[B]:GLU:HG3	8:A:32:HOH:O	2.18	0.44

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Atom-1	Atom-2	Interatomic	Clash	
Atom-1	Atom-2	${ m distance}({ m \AA})$	overlap (Å)	
1:A:257:LEU:O	1:A:258:LEU:CB	2.66	0.44	
1:A:200:LYS:HE2	8:A:336:HOH:O	2.18	0.43	
1:A:184:ILE:CG2	1:A:252[A]:THR:HG23	2.50	0.42	

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	1100111 1		Clash overlap (Å)
8:A:334:HOH:O	8:A:336:HOH:O[2_546]	1.49	0.71

## 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 Allowe		Outliers	Percentiles	
1	A	212/204 (104%)	208 (98%)	4 (2%)	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 side chain 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	187/177 (106%)	184 (98%)	3 (2%)	62 41	

All (3) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	A	61	ARG
1	A	124[A]	PHE
1	A	124[B]	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	144	HIS
1	A	231	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 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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	l Type Chain Res		Res Link Bond lengths			Bond angles								
MIOI	Type	Chain	nes	ries	ries	ries	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z  > 2
6	GOL	A	2	-	5,5,5	0.28	0	5,5,5	0.40	0				
6	GOL	A	260	-	5,5,5	0.54	0	5,5,5	0.54	0				
7	XL3	В	1	-	3,3,3	0.45	0	2,2,2	1.86	1 (50%)				
5	AKG	A	259	4	9,9,9	1.95	1 (11%)	11,11,11	1.40	1 (9%)				



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	2	-	-	0/4/4/4	-
6	GOL	A	260	-	-	0/4/4/4	-
7	XL3	В	1	-	-	0/1/1/1	-
5	AKG	A	259	4	-	1/9/9/9	=

#### All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\operatorname{Observed}(\text{\AA})$	Ideal(A)
5	A	259	AKG	C2-C1	-5.68	1.46	1.53

#### All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
7	В	1	XL3	CB-CC-SG	-2.62	111.01	113.74
5	A	259	AKG	O2-C1-C2	2.61	121.12	113.97

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	259	AKG	C3-C4-C5-O3

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,  $95^{th}$  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>	# RSRZ > 2	$OWAB(A^2)$	Q < 0.9
1	A	204/204 (100%)	0.12	8 (3%) 39 36	12, 21, 37, 44	0
2	В	14/14 (100%)	0.18	0 100 100	16, 28, 38, 40	0
3	С	14/14 (100%)	-0.46	0 100 100	21, 27, 32, 33	0
All	All	232/232 (100%)	0.09	8 (3%) 45 42	12, 23, 38, 44	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	124[A]	PHE	9.4
1	A	258	LEU	7.1
1	A	125	SER	4.5
1	A	126	GLY	4.4
1	A	257	LEU	4.1
1	A	127	LEU	4.0
1	A	210	ARG	3.2
1	A	168	ILE	2.3

## 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	$\operatorname{Res}$	Atoms	RSCC	RSR	${f B-factors(\AA^2)}$	Q<0.9
6	GOL	A	2	6/6	0.81	0.14	43,45,49,50	0
6	GOL	A	260	6/6	0.97	0.07	18,21,21,22	0
5	AKG	A	259	10/10	0.97	0.11	15,16,17,19	0
7	XL3	В	1	4/4	0.99	0.11	16,17,17,17	0
4	MN	A	1	1/1	1.00	0.07	16,16,16,16	0

## 6.5 Other polymers (i)

There are no such residues in this entry.

