



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 11, 2021 – 04:31 PM EDT

PDB ID : 2ILM  
Title : Factor Inhibiting HIF-1 Alpha D201A Mutant in Complex with FE(II), Alpha-Ketoglutarate and HIF-1 Alpha 35mer  
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Deposited on : 2006-10-03  
Resolution : 2.30 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](mailto: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.

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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 : 2.23.2  
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) : 2.23.2

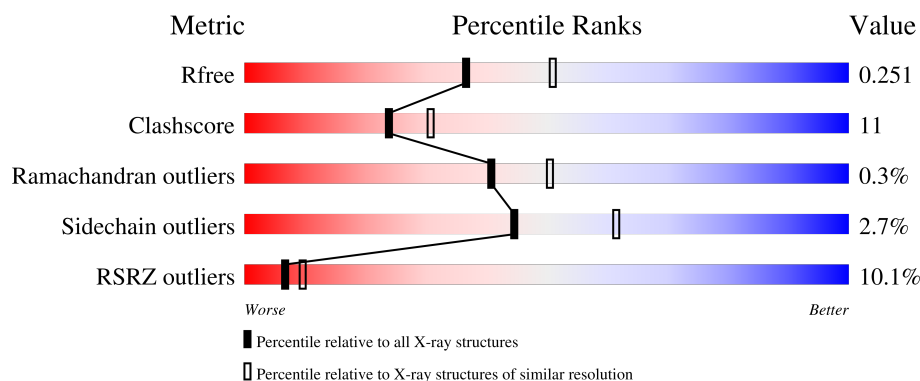
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	349	<div> <div>9%</div> <div>74%</div> <div>23%</div> <div>..</div> </div>
2	S	41	<div> <div>5%</div> <div>10%</div> <div>5%</div> <div>85%</div> </div>

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 2910 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 Hypoxia-inducible factor 1 alpha inhibitor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	0	0	0
			2653	1707	447	488	11			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	201	ALA	ASP	engineered mutation	UNP Q9NWT6

- Molecule 2 is a protein called Hypoxia-inducible factor 1 alpha.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	S	6	Total	C	N	O	0	0	0
			39	23	6	10			

- Molecule 3 is FE (II) ION (three-letter code: FE2) (formula: Fe).

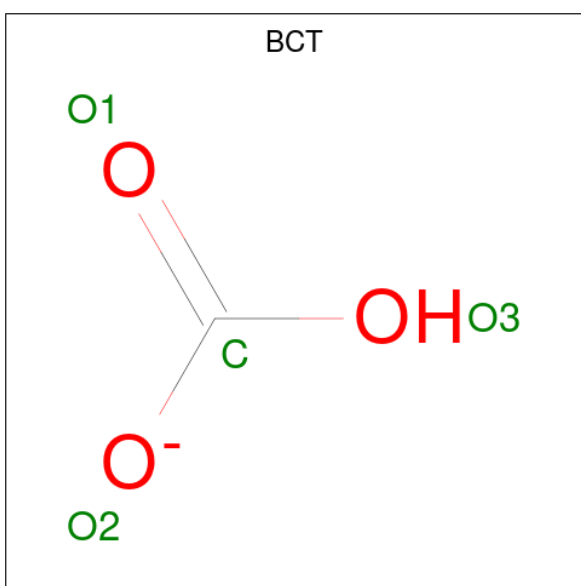
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Fe	0	0
			1	1		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- Molecule 5 is BICARBONATE ION (three-letter code: BCT) (formula:  $\text{CHO}_3$ ).



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

- Molecule 6 is 2-OXOGLUTARIC ACID (three-letter code: AKG) (formula:  $\text{C}_5\text{H}_6\text{O}_5$ ).



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

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



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

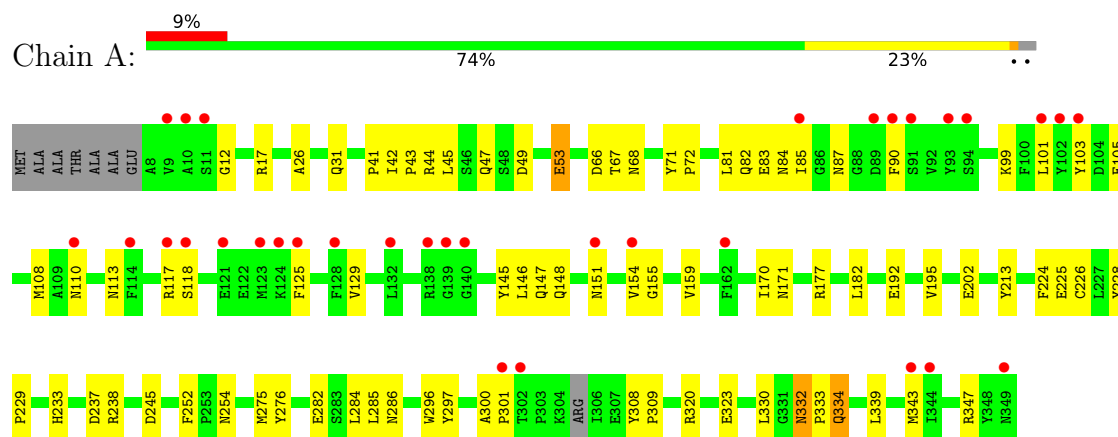
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	172	Total 172	O 172	0	0
8	S	3	Total 3	O 3	0	0

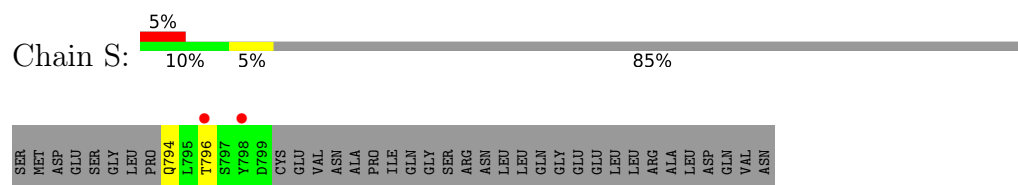
### 3 Residue-property plots [i](#)

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: Hypoxia-inducible factor 1 alpha inhibitor



- Molecule 2: Hypoxia-inducible factor 1 alpha



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	86.53Å 86.53Å 147.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.27 – 2.30 47.11 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.8 (43.27-2.30) 95.0 (47.11-2.30)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.03	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.20 (at 2.29Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.209 , 0.267 0.195 , 0.251	Depositor DCC
$R_{free}$ test set	2529 reflections (9.88%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	47.4	Xtriage
Anisotropy	0.271	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 78.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	2910	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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: BCT, GOL, FE2, SO4, 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).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.47	0/2733	0.66	2/3715 (0.1%)
2	S	0.42	0/38	0.57	0/51
All	All	0.47	0/2771	0.66	2/3766 (0.1%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	347	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	A	347	ARG	NE-CZ-NH1	5.63	123.11	120.30

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	2653	0	2433	57	0
2	S	39	0	30	1	0
3	A	1	0	0	0	0
4	A	10	0	0	0	0
5	A	4	0	0	0	0
6	A	10	0	4	0	0
7	A	18	0	21	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	172	0	0	4	0
8	S	3	0	0	0	0
All	All	2910	0	2488	58	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 11.

All (58) 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:53:GLU:HG3	1:A:177:ARG:HH22	1.37	0.87
1:A:320:ARG:CG	1:A:320:ARG:NE	2.47	0.78
1:A:105:GLU:HA	1:A:108:MET:HG3	1.64	0.77
1:A:332:ASN:HD22	1:A:333:PRO:HD2	1.52	0.73
1:A:334:GLN:HE21	1:A:334:GLN:HA	1.55	0.71
1:A:155:GLY:O	1:A:159:VAL:HG23	1.93	0.68
1:A:125:PHE:O	1:A:129:VAL:HG23	2.00	0.61
1:A:53:GLU:HG3	1:A:177:ARG:NH2	2.12	0.61
1:A:309:PRO:HA	8:A:1138:HOH:O	2.04	0.58
1:A:332:ASN:HD22	1:A:333:PRO:CD	2.17	0.58
1:A:228:TYR:CE2	1:A:252:PHE:HB3	2.39	0.58
1:A:146:LEU:C	1:A:146:LEU:HD23	2.27	0.55
1:A:26:ALA:HB1	1:A:213:TYR:CZ	2.43	0.53
1:A:71:TYR:HB3	1:A:72:PRO:HD3	1.91	0.53
1:A:81:LEU:C	1:A:83:GLU:H	2.11	0.53
1:A:84:ASN:O	1:A:155:GLY:HA3	2.09	0.53
1:A:53:GLU:CG	1:A:177:ARG:HH22	2.16	0.53
1:A:146:LEU:HD23	1:A:147:GLN:N	2.23	0.53
1:A:45:LEU:HD22	1:A:49:ASP:OD2	2.10	0.52
1:A:275:MET:O	1:A:276:TYR:HB2	2.10	0.51
1:A:226:CYS:HB3	1:A:254:ASN:O	2.12	0.50
1:A:147:GLN:NE2	8:A:1140:HOH:O	2.43	0.50
1:A:202:GLU:C	1:A:275:MET:HG3	2.32	0.50
1:A:228:TYR:CD2	1:A:252:PHE:HB3	2.48	0.49
1:A:99:LYS:HD3	8:A:1061:HOH:O	2.12	0.48
1:A:148:GLN:HE22	7:A:2003:GOL:H2	1.78	0.48
1:A:334:GLN:HA	1:A:334:GLN:NE2	2.25	0.47
1:A:82:GLN:HG2	1:A:82:GLN:O	2.14	0.47
1:A:192:GLU:OE1	1:A:286:ASN:HA	2.14	0.47
1:A:237:ASP:O	1:A:238:ARG:HB2	2.15	0.47
1:A:308:TYR:HA	1:A:309:PRO:C	2.36	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:ALA:HB1	1:A:301:PRO:HD2	1.98	0.45
2:S:794:GLN:C	2:S:796:THR:H	2.19	0.45
1:A:233:HIS:HE1	1:A:323:GLU:O	2.00	0.45
1:A:12:GLY:O	1:A:41:PRO:HB3	2.17	0.44
1:A:31:GLN:HB3	1:A:284:LEU:HB3	1.98	0.44
1:A:66:ASP:N	1:A:66:ASP:OD1	2.50	0.44
1:A:225:GLU:OE1	1:A:225:GLU:N	2.42	0.43
1:A:224:PHE:CE2	1:A:229:PRO:HG3	2.54	0.43
1:A:17:ARG:HD3	1:A:44:ARG:CZ	2.49	0.43
1:A:330:LEU:HD21	1:A:339:LEU:HD13	2.01	0.43
1:A:182:LEU:HD13	1:A:297:TYR:CZ	2.54	0.42
1:A:85:ILE:HG12	1:A:90:PHE:HZ	1.85	0.42
1:A:103:TYR:HE2	1:A:117:ARG:H	1.66	0.42
1:A:31:GLN:HG2	1:A:285:LEU:O	2.20	0.42
1:A:182:LEU:HD13	1:A:297:TYR:CE1	2.54	0.42
1:A:101:LEU:HD23	1:A:103:TYR:HE1	1.85	0.42
1:A:67:THR:O	1:A:68:ASN:HB2	2.20	0.42
1:A:334:GLN:HE21	1:A:334:GLN:CA	2.20	0.41
1:A:110:ASN:N	1:A:110:ASN:ND2	2.69	0.41
1:A:85:ILE:O	1:A:154:VAL:HA	2.21	0.41
1:A:195:VAL:HG22	1:A:282:GLU:HB3	2.04	0.40
1:A:339:LEU:HD23	1:A:343:MET:HE2	2.02	0.40
1:A:87:ASN:HA	8:A:1160:HOH:O	2.21	0.40
1:A:99:LYS:NZ	1:A:245:ASP:OD2	2.53	0.40
1:A:103:TYR:HE2	1:A:118:SER:H	1.69	0.40
1:A:170:ILE:CG2	1:A:171:ASN:N	2.85	0.40
1:A:42:ILE:HA	1:A:43:PRO:HD3	1.95	0.40

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	A	337/349 (97%)	314 (93%)	22 (6%)	1 (0%)	41	50
2	S	4/41 (10%)	2 (50%)	2 (50%)	0	100	100
All	All	341/390 (87%)	316 (93%)	24 (7%)	1 (0%)	41	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	113	ASN

### 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	A	260/303 (86%)	253 (97%)	7 (3%)	44	61
2	S	4/36 (11%)	4 (100%)	0	100	100
All	All	264/339 (78%)	257 (97%)	7 (3%)	44	61

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

Mol	Chain	Res	Type
1	A	47	GLN
1	A	53	GLU
1	A	145	TYR
1	A	151	ASN
1	A	296	TRP
1	A	332	ASN
1	A	334	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	ASN
1	A	110	ASN
1	A	147	GLN

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Mol	Chain	Res	Type
1	A	203	GLN
1	A	332	ASN
1	A	334	GLN

### 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 8 ligands modelled in this entry, 1 is 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
4	SO4	A	1353	-	4,4,4	0.28	0	6,6,6	0.14	0
7	GOL	A	2003	-	5,5,5	2.13	2 (40%)	5,5,5	0.45	0
7	GOL	A	2001	-	5,5,5	2.31	2 (40%)	5,5,5	1.15	0
7	GOL	A	2002	-	5,5,5	2.13	2 (40%)	5,5,5	0.78	0
5	BCT	A	3000	3	0,3,3	-	-	0,3,3	-	-
6	AKG	A	1351	3	3,9,9	4.73	2 (66%)	4,11,11	2.07	1 (25%)
4	SO4	A	1352	-	4,4,4	0.26	0	6,6,6	0.06	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
7	GOL	A	2003	-	-	0/4/4/4	-
7	GOL	A	2001	-	-	0/4/4/4	-
7	GOL	A	2002	-	-	0/4/4/4	-
6	AKG	A	1351	3	-	1/3/9/9	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	1351	AKG	C3-C2	-7.70	1.39	1.51
7	A	2003	GOL	O1-C1	-3.57	1.27	1.42
7	A	2001	GOL	O1-C1	-3.51	1.27	1.42
7	A	2002	GOL	O1-C1	-3.45	1.27	1.42
7	A	2001	GOL	O2-C2	3.02	1.52	1.43
7	A	2002	GOL	O2-C2	2.88	1.52	1.43
6	A	1351	AKG	C4-C3	-2.60	1.39	1.52
7	A	2003	GOL	O2-C2	2.41	1.50	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1351	AKG	C4-C3-C2	4.03	121.85	113.14

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1351	AKG	C2-C3-C4-C5

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	2003	GOL	1	0

## 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

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	341/349 (97%)	0.46	33 (9%) <b>7</b> <b>10</b>	28, 54, 104, 138	0
2	S	6/41 (14%)	2.09	2 (33%) <b>0</b> <b>0</b>	73, 86, 121, 126	0
All	All	347/390 (88%)	0.49	35 (10%) <b>7</b> <b>9</b>	28, 54, 111, 138	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	90	PHE	6.4
1	A	118	SER	5.8
1	A	140	GLY	5.6
2	S	798	TYR	5.1
1	A	128	PHE	4.9
1	A	102	TYR	4.6
1	A	132	LEU	4.6
1	A	103	TYR	4.4
1	A	125	PHE	4.2
1	A	114	PHE	4.2
1	A	117	ARG	4.0
1	A	301	PRO	3.5
1	A	121	GLU	3.4
1	A	139	GLY	3.3
1	A	11	SER	3.0
1	A	123	MET	2.7
1	A	10	ALA	2.7
1	A	94	SER	2.6
2	S	796	THR	2.5
1	A	85	ILE	2.5
1	A	349	ASN	2.5
1	A	343	MET	2.4
1	A	9	VAL	2.4
1	A	89	ASP	2.4

*Continued on next page...*



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Mol	Chain	Res	Type	RSRZ
1	A	91	SER	2.4
1	A	344	ILE	2.3
1	A	101	LEU	2.3
1	A	124	LYS	2.3
1	A	151	ASN	2.2
1	A	110	ASN	2.2
1	A	138	ARG	2.1
1	A	93	TYR	2.1
1	A	162	PHE	2.1
1	A	154	VAL	2.1
1	A	302	THR	2.0

## 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<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	B-factors( $\text{\AA}^2$ )	Q<0.9
7	GOL	A	2002	6/6	0.68	0.12	86,96,110,112	0
7	GOL	A	2001	6/6	0.71	0.16	83,88,95,116	0
7	GOL	A	2003	6/6	0.74	0.24	83,116,119,119	0
4	SO4	A	1352	5/5	0.90	0.25	137,139,142,143	0
5	BCT	A	3000	4/4	0.92	0.10	53,54,69,73	0
4	SO4	A	1353	5/5	0.92	0.30	96,118,129,129	0
6	AKG	A	1351	10/10	0.98	0.17	35,51,57,83	0
3	FE2	A	350	1/1	0.99	0.09	46,46,46,46	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.