



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:04 AM UTC

PDB ID : 3D8C / pdb_00003d8c
Title : Factor inhibiting HIF-1 alpha D201G mutant in complex with ZN(II), alpha-ketoglutarate and HIF-1 alpha 19mer
Authors : McDonough, M.A.; Chowdhury, R.; Schofield, C.J.
Deposited on : 2008-05-23
Resolution : 2.10 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

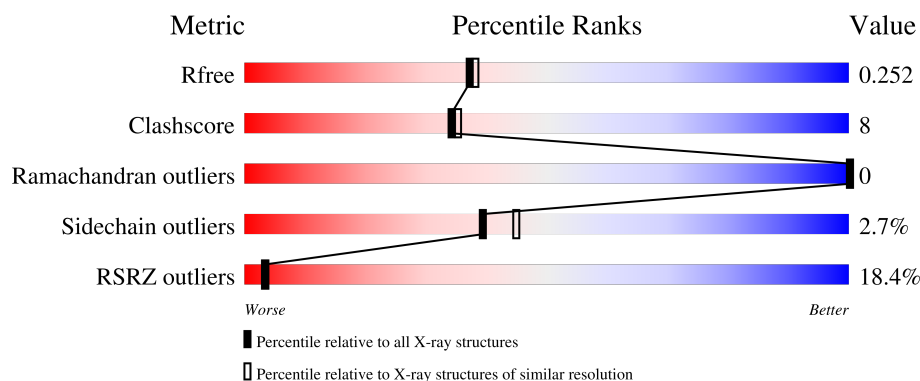
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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 $\leq 5\%$. 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	18% 80% 17% ..
2	B	19	11% 58% 11% 32%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 3072 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
			2725	1752	462	500	11			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q9NWT6
A	2	ALA	-	expression tag	UNP Q9NWT6
A	3	ALA	-	expression tag	UNP Q9NWT6
A	4	THR	-	expression tag	UNP Q9NWT6
A	5	ALA	-	expression tag	UNP Q9NWT6
A	6	ALA	-	expression tag	UNP Q9NWT6
A	7	GLU	-	expression tag	UNP Q9NWT6
A	8	ALA	-	expression tag	UNP Q9NWT6
A	9	VAL	-	expression tag	UNP Q9NWT6
A	10	ALA	-	expression tag	UNP Q9NWT6
A	201	GLY	ASP	engineered mutation	UNP Q9NWT6

- Molecule 2 is a protein called HYPOXIA-INDUCIBLE FACTOR 1 ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	13	Total	C	N	O	S	0	0	0
			97	60	14	22	1			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

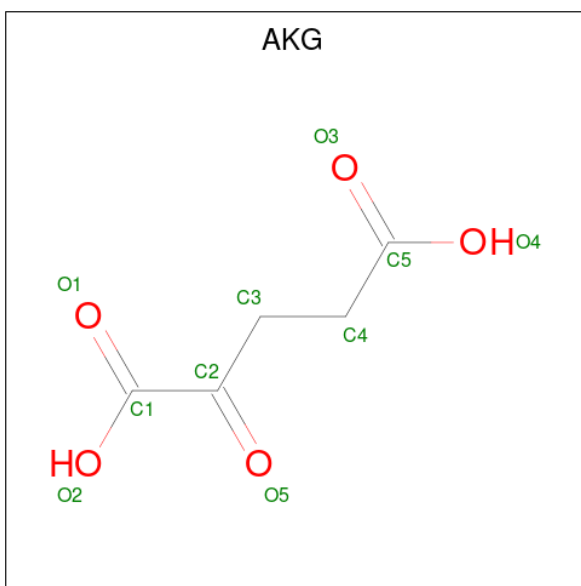
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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 2-OXOGLUTARIC ACID (CCD ID: AKG) (formula: $C_5H_6O_5$).



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

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

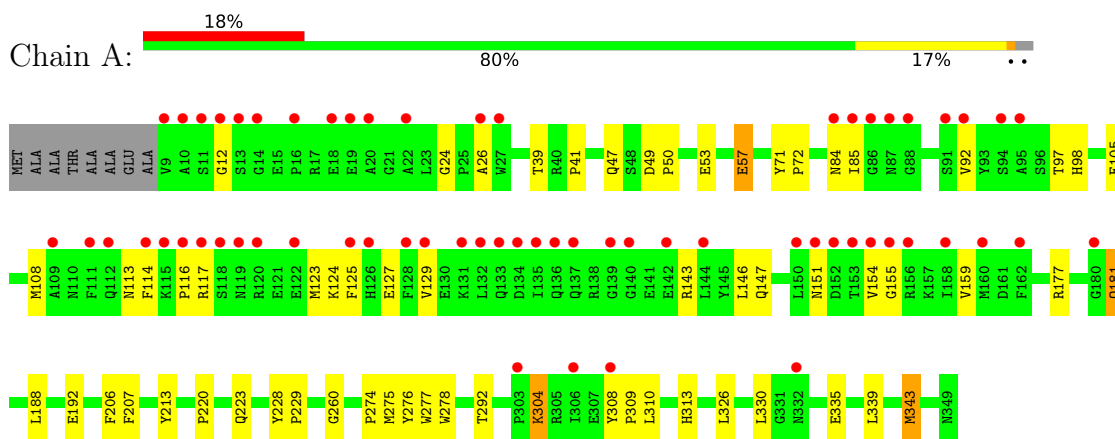
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	214	Total	O	0	0
			214	214		
7	B	9	Total	O	0	0
			9	9		

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, α , β , γ	86.43Å 86.43Å 148.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.65 – 2.10 38.65 – 2.10	Depositor EDS
% Data completeness (in resolution range)	95.2 (38.65-2.10) 95.1 (38.65-2.10)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.218 , 0.256 0.214 , 0.252	Depositor DCC
R_{free} test set	3861 reflections (10.01%)	wwPDB-VP
Wilson B-factor (Å ²)	40.5	Xtriage
Anisotropy	0.351	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 58.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3072	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.57% 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: SO4, AKG, GOL, ZN

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	0.61	1/2807 (0.0%)	0.88	1/3817 (0.0%)
2	B	0.48	0/98	0.80	0/133
All	All	0.61	1/2905 (0.0%)	0.88	1/3950 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	343	MET	SD-CE	-6.41	1.63	1.79

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	206	PHE	N-CA-C	-5.43	99.67	108.52

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	2725	0	2560	41	0
2	B	97	0	86	1	0
3	A	1	0	0	0	0
4	A	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	10	0	4	0	0
6	A	6	0	7	0	0
7	A	214	0	0	5	0
7	B	9	0	0	0	0
All	All	3072	0	2657	42	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 8.

All (42) 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:304:LYS:HD2	1:A:304:LYS:H	1.22	1.00
1:A:57:GLU:HG2	1:A:177:ARG:NH1	2.02	0.73
1:A:181:GLN:HE21	1:A:181:GLN:H	1.39	0.70
1:A:304:LYS:H	1:A:304:LYS:CD	2.00	0.62
1:A:275:MET:O	1:A:276:TYR:HB2	2.06	0.56
1:A:155:GLY:O	1:A:159:VAL:HG23	2.06	0.55
1:A:181:GLN:HE21	1:A:181:GLN:N	2.04	0.54
1:A:71:TYR:HB3	1:A:72:PRO:HD3	1.90	0.54
1:A:147:GLN:HG2	1:A:188:LEU:HD23	1.91	0.53
1:A:57:GLU:HG2	1:A:177:ARG:HH11	1.74	0.52
1:A:220:PRO:HD2	1:A:223:GLN:HG3	1.91	0.52
1:A:339:LEU:CD2	1:A:343:MET:HE2	2.40	0.52
1:A:274:PRO:HB2	1:A:277:TRP:CD1	2.45	0.52
1:A:229:PRO:HB3	1:A:278:TRP:HH2	1.75	0.51
1:A:339:LEU:HD23	1:A:343:MET:HE2	1.93	0.51
1:A:105:GLU:HA	1:A:108:MET:HG3	1.93	0.51
1:A:304:LYS:HD2	1:A:304:LYS:N	2.06	0.51
1:A:97:THR:HB	1:A:117:ARG:HH11	1.76	0.49
1:A:85:ILE:O	1:A:154:VAL:HA	2.12	0.49
1:A:26:ALA:HB1	1:A:213:TYR:CZ	2.49	0.48
1:A:228:TYR:HA	7:A:1016:HOH:O	2.13	0.48
1:A:181:GLN:HB2	7:A:1129:HOH:O	2.12	0.47
1:A:308:TYR:HA	1:A:309:PRO:C	2.40	0.47
1:A:207:PHE:CZ	1:A:292:THR:HG21	2.50	0.47
1:A:12:GLY:O	1:A:41:PRO:HB3	2.15	0.46
1:A:146:LEU:C	1:A:146:LEU:HD23	2.41	0.46
1:A:276:TYR:CD2	1:A:313:HIS:HB2	2.51	0.46
1:A:114:PHE:CE2	1:A:116:PRO:HG3	2.51	0.45
1:A:53:GLU:HG3	1:A:177:ARG:HH22	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:GLU:HG2	7:A:1073:HOH:O	2.17	0.43
1:A:124:LYS:O	1:A:127:GLU:HB2	2.19	0.43
1:A:330:LEU:HD13	1:A:335:GLU:HB3	2.01	0.42
1:A:92:VAL:HB	1:A:123:MET:HE3	2.00	0.42
2:B:795:LEU:HD12	2:B:795:LEU:HA	1.90	0.42
1:A:125:PHE:O	1:A:129:VAL:HG23	2.20	0.42
1:A:39:THR:HA	1:A:260:GLY:O	2.20	0.42
1:A:98:HIS:HA	1:A:143:ARG:CD	2.50	0.42
1:A:84:ASN:O	1:A:155:GLY:HA3	2.20	0.41
1:A:98:HIS:HA	1:A:143:ARG:HD3	2.03	0.41
1:A:24:GLY:N	7:A:1141:HOH:O	2.54	0.41
1:A:113:ASN:HB2	7:A:1099:HOH:O	2.21	0.40
1:A:49:ASP:HA	1:A:50:PRO:HD3	1.97	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	339/349 (97%)	329 (97%)	10 (3%)	0	100	100
2	B	11/19 (58%)	8 (73%)	3 (27%)	0	100	100
All	All	350/368 (95%)	337 (96%)	13 (4%)	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	280/303 (92%)	273 (98%)	7 (2%)	42	48
2	B	11/17 (65%)	10 (91%)	1 (9%)	9	6
All	All	291/320 (91%)	283 (97%)	8 (3%)	39	45

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	57	GLU
1	A	151	ASN
1	A	181	GLN
1	A	304	LYS
1	A	310	LEU
1	A	326	LEU
2	B	796	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	110	ASN
1	A	133	GLN
1	A	147	GLN
1	A	151	ASN
1	A	173	GLN
1	A	181	GLN
1	A	254	ASN
1	A	334	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	GOL	A	2000	-	5,5,5	1.28	1 (20%)	5,5,5	0.39	0
5	AKG	A	1351	3	9,9,9	3.71	4 (44%)	11,11,11	2.01	5 (45%)
4	SO4	A	501	-	4,4,4	0.37	0	6,6,6	0.14	0
4	SO4	A	502	-	4,4,4	0.37	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
5	AKG	A	1351	3	-	1/9/9/9	-
6	GOL	A	2000	-	-	0/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1351	AKG	C3-C2	-9.85	1.40	1.51
5	A	1351	AKG	C4-C3	-3.61	1.39	1.51
5	A	1351	AKG	O2-C1	-2.59	1.23	1.30
5	A	1351	AKG	O4-C5	-2.25	1.23	1.30
6	A	2000	GOL	O3-C3	-2.15	1.33	1.42

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1351	AKG	C4-C3-C2	3.87	120.17	112.91
5	A	1351	AKG	C3-C2-C1	2.63	120.32	115.86
5	A	1351	AKG	O4-C5-C4	2.54	122.04	114.00
5	A	1351	AKG	O2-C1-C2	2.24	119.79	113.59
5	A	1351	AKG	O4-C5-O3	-2.20	117.67	123.33

There are no chirality outliers.

All (1) torsion outliers are listed below:

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

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

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, 95th 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 ²)	Q<0.9
1	A	341/349 (97%)	0.86	63 (18%) 3 3	30, 50, 99, 115	0
2	B	13/19 (68%)	0.93	2 (15%) 5 5	44, 58, 70, 71	0
All	All	354/368 (96%)	0.87	65 (18%) 3 3	30, 50, 99, 115	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	12	GLY	5.3
1	A	112	GLN	5.3
1	A	9	VAL	4.7
1	A	116	PRO	4.7
1	A	152	ASP	4.5
1	A	135	ILE	4.4
1	A	16	PRO	4.3
1	A	134	ASP	4.2
1	A	13	SER	4.1
1	A	114	PHE	3.8
1	A	10	ALA	3.5
1	A	132	LEU	3.5
1	A	136	GLN	3.4
1	A	95	ALA	3.4
1	A	125	PHE	3.4
1	A	118	SER	3.3
1	A	85	ILE	3.3
1	A	115	LYS	3.2
1	A	131	LYS	3.2
1	A	26	ALA	3.2
1	A	153	THR	3.1
1	A	111	PHE	3.1
1	A	92	VAL	3.0
1	A	27	TRP	3.0

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Mol	Chain	Res	Type	RSRZ
2	B	806	ILE	3.0
1	A	129	VAL	2.9
1	A	84	ASN	2.9
1	A	122	GLU	2.7
1	A	151	ASN	2.7
1	A	162	PHE	2.7
1	A	91	SER	2.7
1	A	150	LEU	2.6
1	A	332	ASN	2.6
1	A	155	GLY	2.5
1	A	126	HIS	2.5
1	A	14	GLY	2.5
1	A	306	ILE	2.5
1	A	128	PHE	2.5
1	A	308	TYR	2.4
1	A	94	SER	2.4
1	A	139	GLY	2.4
1	A	18	GLU	2.4
1	A	137	GLN	2.4
1	A	87	ASN	2.4
1	A	119	ASN	2.3
1	A	11	SER	2.3
1	A	144	LEU	2.3
1	A	154	VAL	2.3
1	A	133	GLN	2.3
1	A	303	PRO	2.3
1	A	180	GLY	2.3
1	A	19	GLU	2.3
1	A	142	GLU	2.2
1	A	120	ARG	2.2
1	A	86	GLY	2.1
1	A	156	ARG	2.1
1	A	117	ARG	2.1
1	A	158	ILE	2.1
1	A	109	ALA	2.1
2	B	794	GLN	2.1
1	A	20	ALA	2.1
1	A	88	GLY	2.0
1	A	160	MET	2.0
1	A	22	ALA	2.0
1	A	140	GLY	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 oligosaccharides 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, 95th 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
4	SO4	A	501	5/5	0.89	0.09	95,95,96,97	0
4	SO4	A	502	5/5	0.90	0.11	95,96,96,97	0
6	GOL	A	2000	6/6	0.91	0.23	88,89,91,91	0
5	AKG	A	1351	10/10	0.95	0.08	43,45,48,48	0
3	ZN	A	350	1/1	0.99	0.02	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.