



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:29 AM UTC

PDB ID : 2WA4 / pdb_00002wa4
Title : FACTOR INHIBITING HIF-1 ALPHA WITH N,3-dihydroxybenzamide
Authors : Conejo-Garcia, A.; Lienard, B.M.R.; Clifton, I.J.; McDonough, M.A.; Schofield, C.J.
Deposited on : 2009-02-02
Resolution : 2.50 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

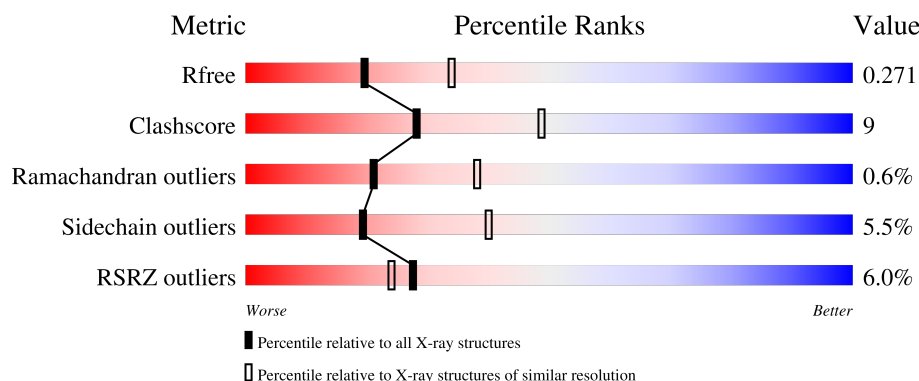
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

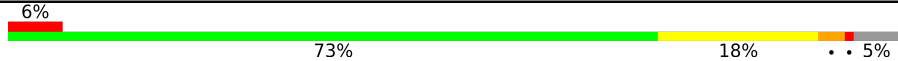
The reported resolution of this entry is 2.50 Å.

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	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 $\leq 5\%$. 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	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2813 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.

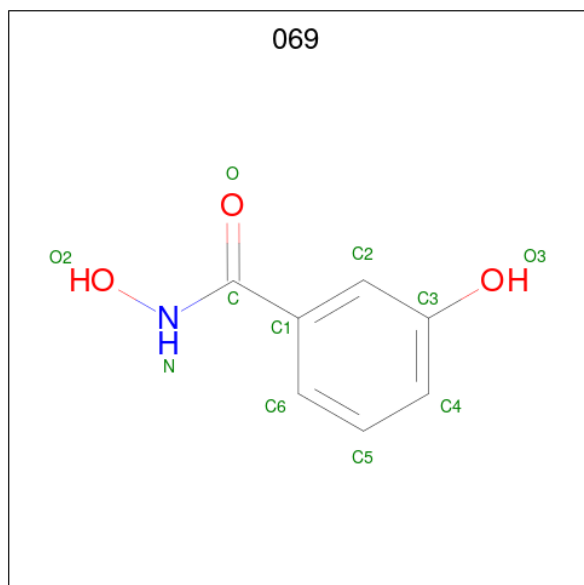
- Molecule 1 is a protein called HYPOXIA-INDUCIBLE FACTOR 1-ALPHA INHIBITOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	332	Total	C	N	O	S	0	1	0
			2731	1752	461	507	11			

- Molecule 2 is FE (II) ION (CCD ID: FE2) (formula: Fe).

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

- Molecule 3 is N,3-DIHYDROXYBENZAMIDE (CCD ID: 069) (formula: C₇H₇NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			11	7	1	3		
3	A	1	Total	C	N	O	0	0
			11	7	1	3		

- 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		

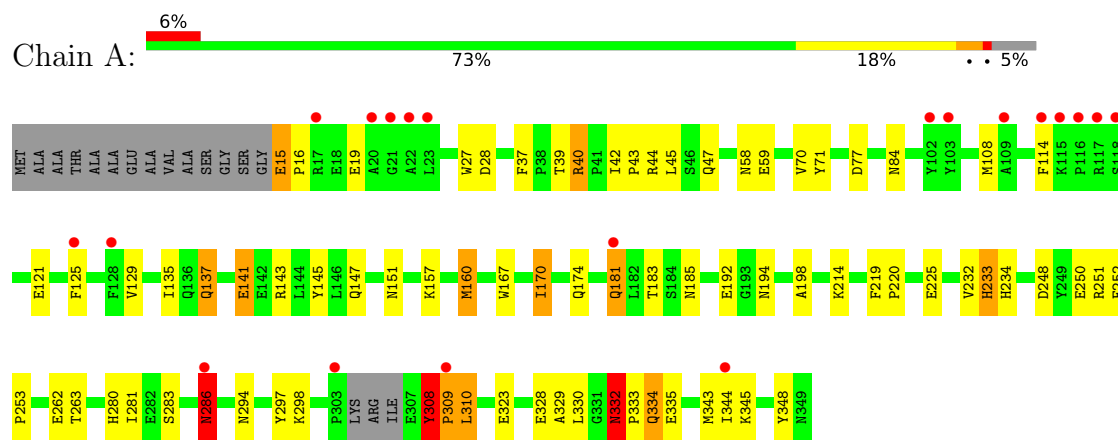
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	54	Total	O	0	0
			54	54		

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



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	87.11Å 87.11Å 150.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	75.38 – 2.50 75.38 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (75.38-2.50) 99.9 (75.38-2.50)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.211 , 0.267 0.212 , 0.271	Depositor DCC
R_{free} test set	1039 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	68.2	Xtriage
Anisotropy	0.244	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 65.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2813	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FE2, 069, SO4

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	1.24	9/2809 (0.3%)	1.16	16/3809 (0.4%)

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 mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	160	MET	SD-CE	9.44	2.03	1.79
1	A	147	GLN	CD-OE1	6.26	1.35	1.23
1	A	334	GLN	CD-OE1	6.17	1.35	1.23
1	A	294	ASN	CG-OD1	5.93	1.34	1.23
1	A	198	ALA	CA-CB	5.92	1.63	1.53
1	A	137	GLN	CD-OE1	5.40	1.33	1.23
1	A	233	HIS	ND1-CE1	5.26	1.37	1.32
1	A	151	ASN	CG-OD1	5.16	1.33	1.23
1	A	294	ASN	CG-ND2	-5.05	1.22	1.33

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	332	ASN	CA-C-N	7.21	127.84	119.47
1	A	332	ASN	C-N-CA	7.21	127.84	119.47
1	A	233	HIS	CB-CG-CD2	-6.34	122.96	131.20
1	A	234	HIS	CA-C-N	-6.23	113.20	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	234	HIS	C-N-CA	-6.23	113.20	119.56
1	A	219	PHE	CA-C-N	-6.07	114.13	120.38
1	A	219	PHE	C-N-CA	-6.07	114.13	120.38
1	A	135	ILE	CB-CA-C	-5.63	104.67	112.04
1	A	286	ASN	N-CA-C	5.35	119.01	111.52
1	A	220	PRO	CA-C-N	5.30	125.61	119.47
1	A	220	PRO	C-N-CA	5.30	125.61	119.47
1	A	15	GLU	CA-C-N	5.29	126.46	119.84
1	A	15	GLU	C-N-CA	5.29	126.46	119.84
1	A	37	PHE	CA-C-N	5.26	125.71	120.14
1	A	37	PHE	C-N-CA	5.26	125.71	120.14
1	A	233	HIS	CB-CG-ND1	5.23	130.54	122.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	308	TYR	Peptide
1	A	309	PRO	Peptide

5.2 Too-close contacts ⓘ

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	2731	0	2602	48	1
2	A	1	0	0	0	0
3	A	22	0	10	3	0
4	A	5	0	0	0	0
5	A	54	0	0	3	0
All	All	2813	0	2612	48	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 9.

All (48) 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:160:MET:CE	1:A:160:MET:SD	2.03	1.47
1:A:344[B]:ILE:CA	1:A:345:LYS:N	1.96	1.28
1:A:343:MET:C	1:A:344[A]:ILE:CA	2.06	1.27
1:A:343:MET:C	1:A:344[B]:ILE:CA	2.33	1.02
1:A:344[A]:ILE:CA	1:A:345:LYS:N	2.50	0.73
1:A:344[B]:ILE:CB	1:A:345:LYS:N	2.50	0.73
1:A:344[B]:ILE:HG23	1:A:348:TYR:HB2	1.71	0.70
1:A:308:TYR:HB3	1:A:309:PRO:CA	2.23	0.68
1:A:141:GLU:OE2	1:A:141:GLU:HA	1.93	0.67
1:A:343:MET:O	1:A:344[B]:ILE:CA	2.44	0.66
1:A:308:TYR:HB3	1:A:309:PRO:HA	1.78	0.65
1:A:233:HIS:HE1	5:A:2047:HOH:O	1.81	0.64
1:A:214:LYS:NZ	1:A:283:SER:OG	2.39	0.56
1:A:39:THR:O	1:A:40:ARG:HD2	2.05	0.56
1:A:108:MET:HE3	1:A:114:PHE:HD2	1.70	0.56
1:A:70:VAL:HG22	5:A:2018:HOH:O	2.06	0.55
1:A:343:MET:O	1:A:344[A]:ILE:CA	2.53	0.55
1:A:167:TRP:HA	1:A:170:ILE:HG22	1.91	0.53
1:A:192:GLU:CD	1:A:286:ASN:HA	2.33	0.53
1:A:214:LYS:CE	3:A:400:069:H4	2.40	0.51
1:A:308:TYR:CB	1:A:309:PRO:HA	2.42	0.50
1:A:232:VAL:HG23	1:A:323:GLU:OE1	2.14	0.48
1:A:183:THR:HG21	1:A:298:LYS:HG3	1.96	0.48
1:A:310:LEU:N	5:A:2043:HOH:O	2.31	0.47
1:A:170:ILE:HG12	1:A:297:TYR:OH	2.14	0.47
1:A:214:LYS:HE3	3:A:400:069:H4	1.96	0.47
1:A:108:MET:HE3	1:A:114:PHE:CD2	2.50	0.46
1:A:84:ASN:ND2	1:A:157:LYS:HD2	2.32	0.45
1:A:248:ASP:CG	1:A:251:ARG:HG3	2.42	0.45
1:A:160:MET:CE	1:A:160:MET:CG	2.93	0.44
1:A:143:ARG:HD2	1:A:194:ASN:OD1	2.18	0.44
1:A:185:ASN:OD1	1:A:185:ASN:C	2.61	0.44
1:A:15:GLU:OE1	1:A:15:GLU:N	2.51	0.43
1:A:214:LYS:HE2	3:A:400:069:H4	2.01	0.43
1:A:174:GLN:OE1	1:A:181:GLN:HA	2.18	0.43
1:A:77:ASP:C	1:A:77:ASP:OD1	2.61	0.43
1:A:280:HIS:C	1:A:281:ILE:HG13	2.44	0.43
1:A:330:LEU:HD13	1:A:335:GLU:HB3	2.00	0.42
1:A:44:ARG:O	1:A:45:LEU:HD23	2.18	0.42
1:A:332:ASN:HA	1:A:333:PRO:HD3	1.90	0.42
1:A:16:PRO:HB3	1:A:27:TRP:O	2.19	0.42
1:A:42:ILE:HA	1:A:43:PRO:HD3	1.88	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:PHE:O	1:A:129:VAL:HG23	2.19	0.42
1:A:58:ASN:O	1:A:59:GLU:C	2.62	0.41
1:A:252:PHE:O	1:A:253:PRO:C	2.62	0.41
1:A:344[B]:ILE:HG22	1:A:345:LYS:N	2.36	0.40
1:A:70:VAL:O	1:A:71:TYR:C	2.65	0.40
1:A:328:GLU:O	1:A:329:ALA:C	2.64	0.40

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	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:TYR:OH	1:A:225:GLU:OE1[5_545]	2.18	0.02

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	325/349 (93%)	301 (93%)	22 (7%)	2 (1%)	21 38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	308	TYR

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	292/304 (96%)	276 (94%)	16 (6%)	19	40

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	GLU
1	A	28	ASP
1	A	40	ARG
1	A	121	GLU
1	A	137	GLN
1	A	141	GLU
1	A	145	TYR
1	A	170	ILE
1	A	181	GLN
1	A	250	GLU
1	A	262	GLU
1	A	263	THR
1	A	286	ASN
1	A	310	LEU
1	A	332	ASN
1	A	334	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	GLN
1	A	126	HIS
1	A	203	GLN
1	A	256	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 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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	$\# Z > 2$
3	069	A	401	2	11,11,11	0.87	0	14,14,14	1.94	5 (35%)
4	SO4	A	402	-	4,4,4	0.54	0	6,6,6	0.33	0
3	069	A	400	2	11,11,11	1.35	1 (9%)	14,14,14	1.06	1 (7%)

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
3	069	A	401	2	-	0/6/6/6	0/1/1/1
3	069	A	400	2	-	0/6/6/6	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	400	069	C1-C	3.19	1.57	1.50

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	069	C1-C-N	-3.21	111.30	116.26
3	A	401	069	C5-C6-C1	2.91	123.23	120.36
3	A	401	069	O2-N-C	2.91	127.02	119.73
3	A	401	069	O-C-N	2.80	128.05	122.90
3	A	401	069	C4-C3-C2	2.49	122.92	120.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	400	069	C4-C3-C2	2.27	122.67	120.19

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	400	069	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, 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	332/349 (95%)	0.27	20 (6%) 27 24	23, 71, 121, 172	1 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	21	GLY	4.8
1	A	118	SER	4.0
1	A	303	PRO	3.9
1	A	23	LEU	3.9
1	A	114	PHE	3.6
1	A	116	PRO	3.2
1	A	20	ALA	3.0
1	A	17	ARG	2.9
1	A	115	LYS	2.9
1	A	102	TYR	2.7
1	A	344[A]	ILE	2.6
1	A	109	ALA	2.5
1	A	103	TYR	2.4
1	A	286	ASN	2.3
1	A	128	PHE	2.3
1	A	117	ARG	2.2
1	A	125	PHE	2.2
1	A	181	GLN	2.1
1	A	309	PRO	2.1
1	A	22	ALA	2.1

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	402	5/5	0.80	0.17	58,59,64,66	5
3	069	A	401	11/11	0.88	0.20	36,56,62,63	11
3	069	A	400	11/11	0.92	0.10	50,53,58,65	0
2	FE2	A	350	1/1	0.98	0.04	53,53,53,53	0

6.5 Other polymers [i](#)

There are no such residues in this entry.