



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:45 PM UTC

PDB ID : 5F4X / pdb_00005f4x
Title : Fructose-1,6-bisphosphate aldolase K229M mutant from rabbit muscle
Authors : LowKam, C.; Arthus-Cartier, G.
Deposited on : 2015-12-03
Resolution : 1.84 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

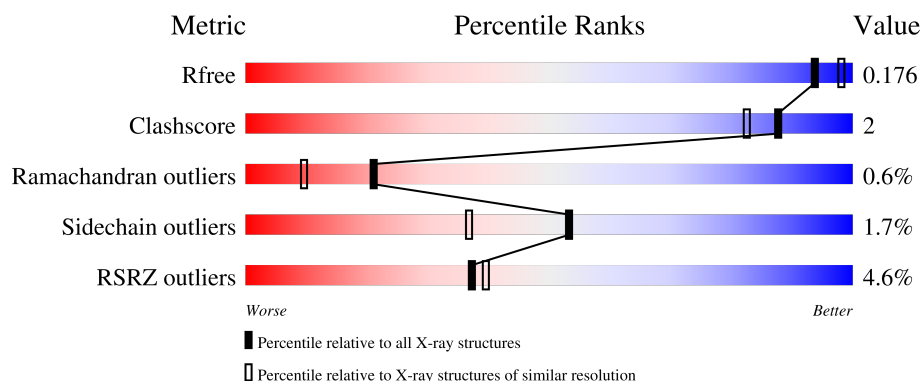
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.84 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	1296 (1.84-1.84)
Clashscore	190562	1329 (1.84-1.84)
Ramachandran outliers	187476	1318 (1.84-1.84)
Sidechain outliers	187428	1318 (1.84-1.84)
RSRZ outliers	180081	1296 (1.84-1.84)

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	363	5% 93% 6%
1	B	363	6% 95% • •
1	C	363	4% 96% •
1	D	363	4% 94% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 24064 atoms, of which 11152 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 Fructose-bisphosphate aldolase A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	363	Total	C	H	N	O	S	0	2	0
			5547	1737	2782	489	526	13			
1	B	363	Total	C	H	N	O	S	0	1	0
			5546	1737	2781	489	526	13			
1	C	363	Total	C	H	N	O	S	0	2	0
			5547	1737	2782	489	526	13			
1	D	363	Total	C	H	N	O	S	0	2	0
			5548	1737	2783	489	526	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	229	MET	LYS	engineered mutation	UNP P00883
B	229	MET	LYS	engineered mutation	UNP P00883
C	229	MET	LYS	engineered mutation	UNP P00883
D	229	MET	LYS	engineered mutation	UNP P00883

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	H	O	0	1
			28	6	16	6		
2	D	1	Total	C	H	O	0	0
			14	3	8	3		

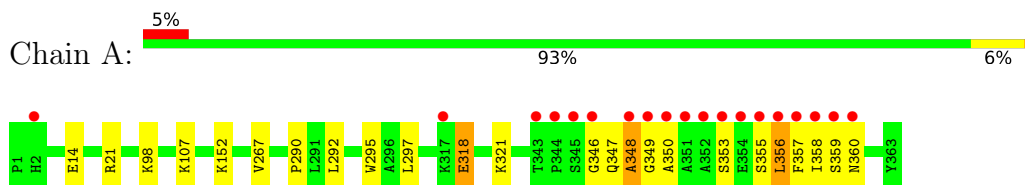
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	481	Total	O	0	0
			481	481		
3	B	470	Total	O	0	0
			470	470		
3	C	474	Total	O	0	0
			474	474		
3	D	409	Total	O	0	0
			409	409		

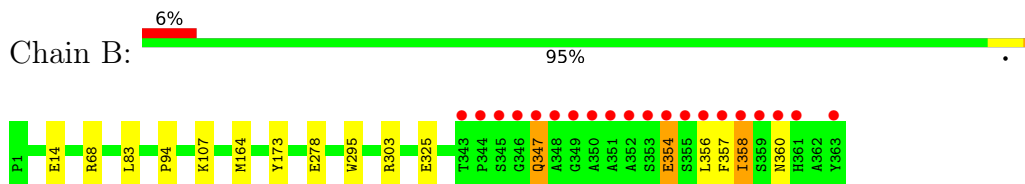
3 Residue-property plots [i](#)

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.

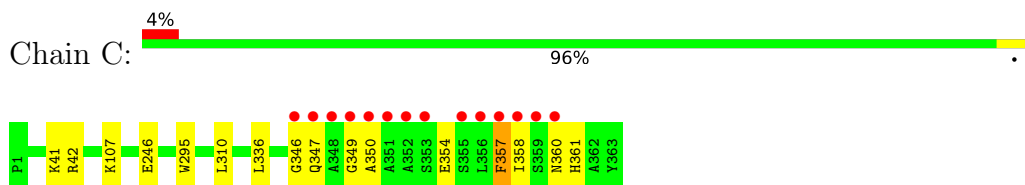
- Molecule 1: Fructose-bisphosphate aldolase A



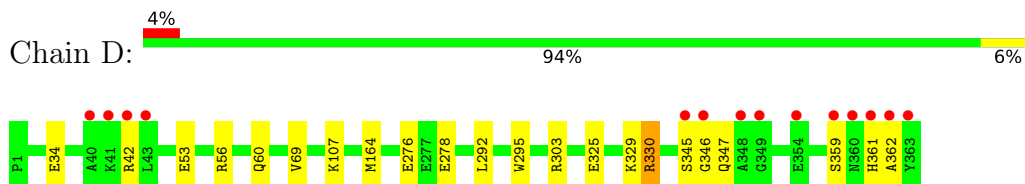
- Molecule 1: Fructose-bisphosphate aldolase A



- Molecule 1: Fructose-bisphosphate aldolase A



- Molecule 1: Fructose-bisphosphate aldolase A



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.48Å 103.19Å 84.53Å 90.00° 98.66° 90.00°	Depositor
Resolution (Å)	35.20 – 1.84 35.20 – 1.84	Depositor EDS
% Data completeness (in resolution range)	99.4 (35.20-1.84) 96.2 (35.20-1.84)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.94 (at 1.84Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.146 , 0.174 0.149 , 0.176	Depositor DCC
R_{free} test set	2002 reflections (1.63%)	wwPDB-VP
Wilson B-factor (Å ²)	17.2	Xtriage
Anisotropy	0.416	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 58.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.016 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	24064	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.33% 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: GOL

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	0/2828	0.73	2/3831 (0.1%)
1	B	0.46	0/2819	0.71	0/3819
1	C	0.46	0/2828	0.71	2/3831 (0.1%)
1	D	0.45	0/2828	0.70	1/3831 (0.0%)
All	All	0.46	0/11303	0.71	5/15312 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	362	ALA	N-CA-C	-6.91	104.59	113.16
1	A	358	ILE	N-CA-C	6.54	116.56	110.42
1	A	353	SER	N-CA-C	-5.20	99.91	108.49
1	C	357	PHE	N-CA-C	5.16	117.82	107.41
1	C	361	HIS	N-CA-C	-5.02	107.20	113.38

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	2765	2782	2775	13	0
1	B	2765	2781	2780	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2765	2782	2775	5	0
1	D	2765	2783	2775	15	0
2	B	12	16	16	0	0
2	D	6	8	8	0	0
3	A	481	0	0	6	1
3	B	470	0	0	8	1
3	C	474	0	0	1	0
3	D	409	0	0	6	0
All	All	12912	11152	11129	45	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 2.

All (45) 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:D:56:ARG:NH1	3:D:501:HOH:O	1.88	1.04
1:D:329:LYS:NZ	3:D:503:HOH:O	2.21	0.72
1:B:278:GLU:OE1	3:B:501:HOH:O	2.08	0.71
1:D:42:ARG:NH1	3:D:506:HOH:O	2.28	0.67
1:A:21:ARG:NH1	3:A:405:HOH:O	2.24	0.67
1:D:60:GLN:OE1	3:D:501:HOH:O	2.13	0.67
1:B:278:GLU:OE2	3:B:502:HOH:O	2.13	0.66
1:A:318:GLU:OE1	3:A:401:HOH:O	2.13	0.65
1:B:164[A]:MET:SD	3:B:819:HOH:O	2.55	0.63
1:B:68:ARG:NH1	3:B:507:HOH:O	2.31	0.59
1:A:14:GLU:OE2	3:A:402:HOH:O	2.17	0.58
1:B:357:PHE:CE2	1:B:358:ILE:HG13	2.40	0.57
1:B:354:GLU:O	3:B:503:HOH:O	2.17	0.57
1:A:152:LYS:NZ	3:A:406:HOH:O	2.24	0.56
1:C:336:LEU:HD21	1:C:347:GLN:OE1	2.05	0.56
1:D:34:GLU:CD	1:D:42:ARG:HH22	2.13	0.56
1:D:346:GLY:HA2	1:D:347:GLN:HB2	1.87	0.55
1:D:53:GLU:OE2	1:D:56:ARG:NH2	2.41	0.54
1:C:246:GLU:OE1	3:C:401:HOH:O	2.19	0.53
1:A:98:LYS:NZ	3:A:409:HOH:O	2.40	0.50
1:B:14:GLU:OE2	3:B:504:HOH:O	2.20	0.50
1:A:360:ASN:N	3:A:411:HOH:O	2.44	0.49
1:B:357:PHE:CD2	1:B:358:ILE:HG13	2.47	0.49
1:A:290:PRO:HG3	1:A:357:PHE:HB2	1.93	0.49
1:D:325:GLU:O	1:D:329:LYS:HG3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:68:ARG:NH2	1:B:325:GLU:OE2	2.41	0.48
1:B:347:GLN:NE2	3:B:513:HOH:O	2.46	0.48
1:D:42:ARG:HD3	1:D:303:ARG:HD3	1.94	0.48
1:D:164[A]:MET:SD	3:D:803:HOH:O	2.61	0.45
1:A:349:GLY:O	1:A:350:ALA:HB3	2.17	0.45
1:A:355:SER:O	1:A:356:LEU:HG	2.15	0.45
1:B:83:LEU:HD12	1:B:94:PRO:HG3	1.99	0.45
1:C:349:GLY:O	1:C:350:ALA:HB3	2.18	0.44
1:A:347:GLN:O	1:A:348:ALA:HB2	2.18	0.44
1:D:346:GLY:CA	1:D:347:GLN:HB2	2.49	0.43
1:A:267:VAL:HB	1:A:297:LEU:HD23	2.00	0.43
1:C:42:ARG:HG3	1:C:310:LEU:CD2	2.49	0.43
1:D:278:GLU:OE1	3:D:502:HOH:O	2.21	0.42
1:D:53:GLU:CD	1:D:56:ARG:HH21	2.27	0.42
1:D:359:SER:HB2	1:D:361:HIS:NE2	2.35	0.42
1:A:290:PRO:HB3	1:A:355:SER:C	2.45	0.41
1:D:276:GLU:HB3	1:D:330:ARG:HD2	2.03	0.41
1:B:303:ARG:HD3	3:B:552:HOH:O	2.20	0.41
1:C:357:PHE:O	1:C:358:ILE:HG13	2.21	0.41
1:A:357:PHE:HE2	1:A:359:SER:HB2	1.86	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)
3:A:506:HOH:O	3:B:549:HOH:O[1_655]	2.15	0.05

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	363/363 (100%)	345 (95%)	15 (4%)	3 (1%)	16 5

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	362/363 (100%)	346 (96%)	14 (4%)	2 (1%)	21	9
1	C	363/363 (100%)	348 (96%)	13 (4%)	2 (1%)	21	9
1	D	363/363 (100%)	353 (97%)	9 (2%)	1 (0%)	36	25
All	All	1451/1452 (100%)	1392 (96%)	51 (4%)	8 (1%)	21	9

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	348	ALA
1	B	347	GLN
1	A	346	GLY
1	B	358	ILE
1	A	356	LEU
1	C	346	GLY
1	C	354	GLU
1	D	345	SER

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	293/291 (101%)	288 (98%)	5 (2%)	53	37
1	B	292/291 (100%)	286 (98%)	6 (2%)	47	30
1	C	293/291 (101%)	289 (99%)	4 (1%)	59	45
1	D	293/291 (101%)	288 (98%)	5 (2%)	53	37
All	All	1171/1164 (101%)	1151 (98%)	20 (2%)	53	37

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	107	LYS
1	A	292	LEU
1	A	295	TRP

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Mol	Chain	Res	Type
1	A	318	GLU
1	A	321	LYS
1	B	107	LYS
1	B	173	TYR
1	B	295	TRP
1	B	354	GLU
1	B	356	LEU
1	B	360	ASN
1	C	41	LYS
1	C	107	LYS
1	C	295	TRP
1	C	360	ASN
1	D	69	VAL
1	D	107	LYS
1	D	292	LEU
1	D	295	TRP
1	D	330	ARG

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

Mol	Chain	Res	Type
1	B	180	ASN
1	B	324	GLN
1	C	136	GLN
1	C	180	ASN
1	D	306	GLN
1	D	360	ASN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

3 ligands are modelled in this entry.

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	$\# Z > 2$
2	GOL	B	401[A]	-	5,5,5	0.43	0	5,5,5	0.73	0
2	GOL	B	401[B]	-	5,5,5	0.37	0	5,5,5	0.20	0
2	GOL	D	401	-	5,5,5	0.46	0	5,5,5	0.61	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	GOL	B	401[A]	-	-	2/4/4/4	-
2	GOL	B	401[B]	-	-	2/4/4/4	-
2	GOL	D	401	-	-	1/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401[B]	GOL	O1-C1-C2-C3
2	B	401[B]	GOL	O1-C1-C2-O2
2	B	401[A]	GOL	C1-C2-C3-O3
2	B	401[A]	GOL	O2-C2-C3-O3
2	D	401	GOL	O2-C2-C3-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

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	363/363 (100%)	-0.32	19 (5%)	33 34	8, 21, 56, 158	9 (2%)
1	B	363/363 (100%)	-0.36	20 (5%)	30 33	10, 20, 65, 184	10 (2%)
1	C	363/363 (100%)	-0.40	14 (3%)	43 46	9, 20, 59, 194	9 (2%)
1	D	363/363 (100%)	-0.23	14 (3%)	43 46	9, 25, 66, 138	3 (0%)
All	All	1452/1452 (100%)	-0.33	67 (4%)	37 39	8, 21, 65, 194	31 (2%)

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	351	ALA	8.2
1	C	348	ALA	7.6
1	C	358	ILE	6.5
1	C	351	ALA	6.3
1	C	346	GLY	6.0
1	A	357	PHE	6.0
1	B	349	GLY	5.9
1	C	347	GLN	5.6
1	C	350	ALA	5.5
1	B	358	ILE	5.5
1	A	355	SER	5.5
1	A	356	LEU	5.4
1	B	356	LEU	5.4
1	C	357	PHE	5.4
1	A	352	ALA	4.9
1	D	346	GLY	4.8
1	D	363	TYR	4.6
1	B	360	ASN	4.6
1	C	349	GLY	4.6
1	B	352	ALA	4.6
1	B	348	ALA	4.5

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Mol	Chain	Res	Type	RSRZ
1	D	362	ALA	4.5
1	C	352	ALA	4.4
1	A	344	PRO	4.4
1	B	350	ALA	4.4
1	A	359	SER	4.3
1	C	353	SER	4.3
1	B	355	SER	4.3
1	A	358	ILE	4.1
1	B	357	PHE	4.1
1	A	345	SER	4.0
1	A	350	ALA	4.0
1	C	356	LEU	4.0
1	B	346	GLY	3.9
1	D	345	SER	3.8
1	A	348	ALA	3.8
1	A	351	ALA	3.8
1	D	361	HIS	3.6
1	B	353	SER	3.4
1	B	359	SER	3.4
1	B	344	PRO	3.3
1	A	360	ASN	3.2
1	A	349	GLY	3.2
1	D	348	ALA	3.1
1	A	353	SER	3.1
1	C	355	SER	3.1
1	A	346	GLY	3.1
1	D	349	GLY	3.1
1	D	41	LYS	3.0
1	D	42	ARG	3.0
1	A	354	GLU	3.0
1	C	359	SER	2.9
1	D	360	ASN	2.9
1	D	354	GLU	2.6
1	A	317	LYS	2.4
1	C	360	ASN	2.3
1	D	359	SER	2.3
1	B	347	GLN	2.3
1	D	43	LEU	2.2
1	B	361	HIS	2.2
1	A	2	HIS	2.1
1	B	343	THR	2.1
1	B	354	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	40	ALA	2.1
1	B	345	SER	2.0
1	B	363	TYR	2.0
1	A	343	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 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
2	GOL	B	401[A]	6/6	0.90	0.10	11,17,23,26	0
2	GOL	B	401[B]	6/6	0.90	0.10	13,16,21,23	14
2	GOL	D	401	6/6	0.93	0.08	14,20,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.