



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 09:27 AM UTC

PDB ID : 3ZGB / pdb_00003zgb
Title : Greater efficiency of photosynthetic carbon fixation due to single amino acid substitution
Authors : Paulus, J.K.; Schlieper, D.; Groth, G.
Deposited on : 2012-12-17
Resolution : 2.71 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

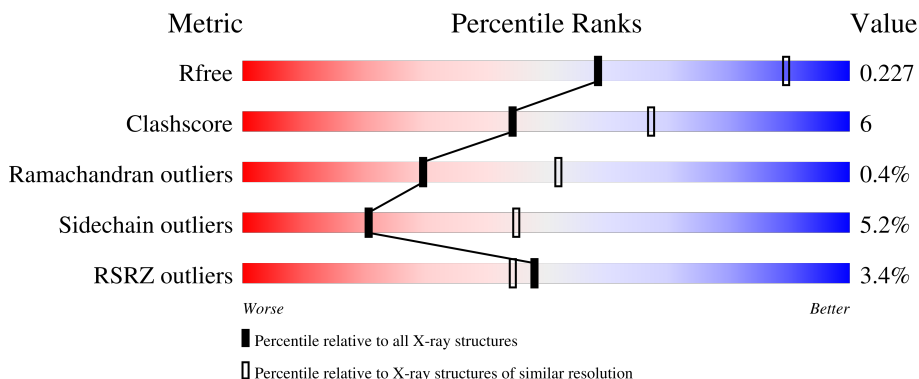
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.71 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	4348 (2.74-2.70)
Clashscore	190562	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
Sidechain outliers	187428	4585 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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	972	
1	B	972	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 14852 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 PHOSPHOENOLPYRUVATE CARBOXYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	920	Total	C	N	O	S	0	0	0
			7396	4691	1289	1382	34			
1	B	915	Total	C	N	O	S	0	0	0
			7354	4663	1282	1375	34			

There are 26 discrepancies between the modelled and reference sequences:

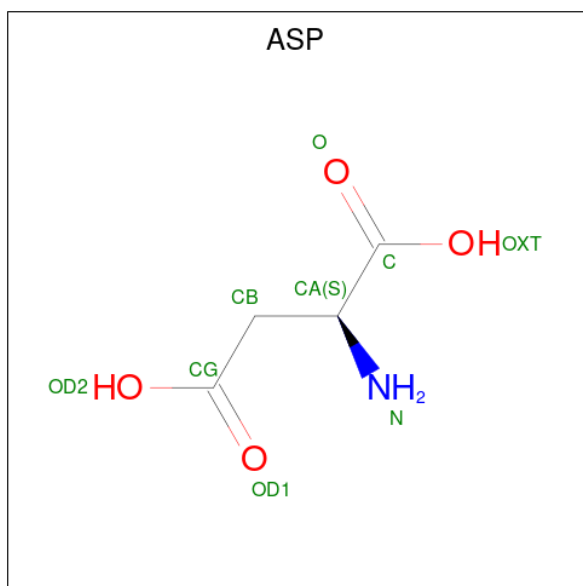
Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	MET	-	expression tag	UNP Q01647
A	-4	GLY	-	expression tag	UNP Q01647
A	-3	HIS	-	expression tag	UNP Q01647
A	-2	HIS	-	expression tag	UNP Q01647
A	-1	HIS	-	expression tag	UNP Q01647
A	0	HIS	-	expression tag	UNP Q01647
A	1	HIS	-	expression tag	UNP Q01647
A	2	HIS	-	expression tag	UNP Q01647
A	3	HIS	-	expression tag	UNP Q01647
A	4	HIS	-	expression tag	UNP Q01647
A	5	HIS	-	expression tag	UNP Q01647
A	?	-	LYS	SEE REMARK 999	UNP Q01647
A	291	ASN	HIS	SEE REMARK 999	UNP Q01647
B	-5	MET	-	expression tag	UNP Q01647
B	-4	GLY	-	expression tag	UNP Q01647
B	-3	HIS	-	expression tag	UNP Q01647
B	-2	HIS	-	expression tag	UNP Q01647
B	-1	HIS	-	expression tag	UNP Q01647
B	0	HIS	-	expression tag	UNP Q01647
B	1	HIS	-	expression tag	UNP Q01647
B	2	HIS	-	expression tag	UNP Q01647
B	3	HIS	-	expression tag	UNP Q01647
B	4	HIS	-	expression tag	UNP Q01647
B	5	HIS	-	expression tag	UNP Q01647
B	?	-	LYS	SEE REMARK 999	UNP Q01647

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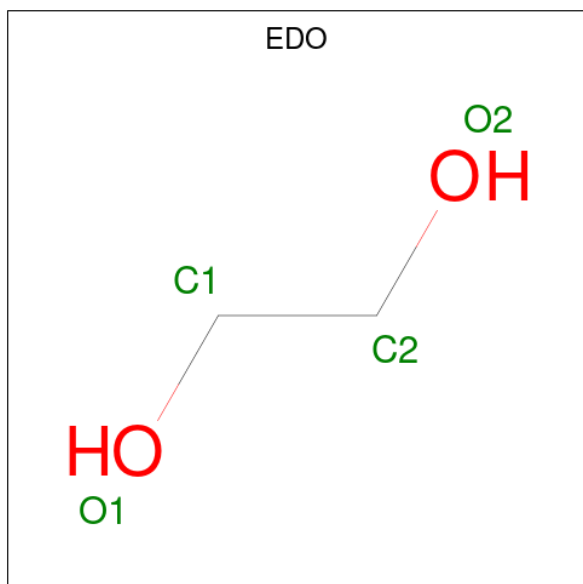
Chain	Residue	Modelled	Actual	Comment	Reference
B	291	ASN	HIS	SEE REMARK 999	UNP Q01647

- Molecule 2 is ASPARTIC ACID (CCD ID: ASP) (formula: $C_4H_7NO_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			9	4	1	4		
2	B	1	Total	C	N	O	0	0
			9	4	1	4		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0

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



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

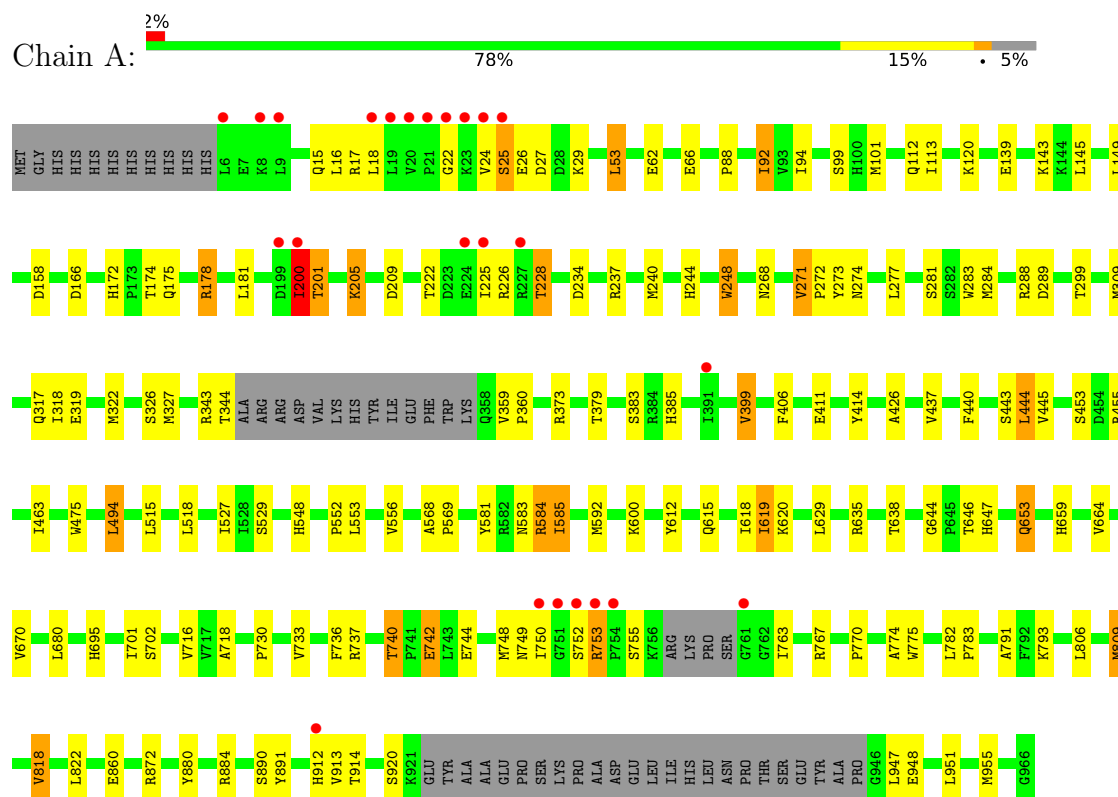
- Molecule 5 is water.

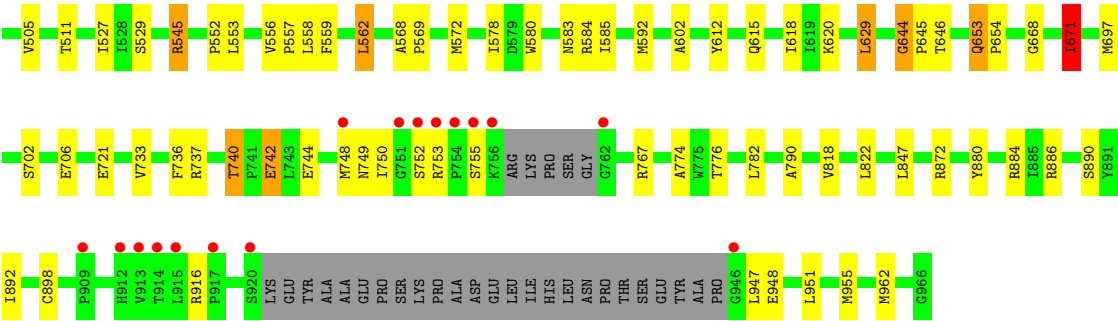
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	35	Total O 35 35	0	0
5	B	21	Total O 21 21	0	0

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: PHOSPHOENOLPYRUVATE CARBOXYLASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	165.55Å 121.70Å 132.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	78.72 – 2.71 78.72 – 2.71	Depositor EDS
% Data completeness (in resolution range)	100.0 (78.72-2.71) 100.0 (78.72-2.71)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	10.98 (at 2.73Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.188 , 0.230 0.188 , 0.227	Depositor DCC
R_{free} test set	1486 reflections (2.03%)	wwPDB-VP
Wilson B-factor (Å ²)	45.2	Xtriage
Anisotropy	0.657	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 39.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14852	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.92% 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: EDO, 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	1.20	19/7549 (0.3%)	1.11	20/10205 (0.2%)
1	B	1.08	4/7506 (0.1%)	1.07	25/10148 (0.2%)
All	All	1.14	23/15055 (0.2%)	1.09	45/20353 (0.2%)

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	B	0	1

The worst 5 of 23 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	659	HIS	CG-CD2	7.57	1.44	1.35
1	A	172	HIS	CG-CD2	7.22	1.43	1.35
1	A	444	LEU	N-CA	7.12	1.55	1.46
1	A	244	HIS	CG-CD2	6.62	1.43	1.35
1	A	475	TRP	NE1-CE2	-6.42	1.30	1.37

The worst 5 of 45 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	443	SER	CA-C-N	-9.60	105.31	121.92
1	A	443	SER	C-N-CA	-9.60	105.31	121.92
1	A	200	ILE	N-CA-C	9.19	122.22	108.80
1	B	443	SER	CA-C-N	-8.54	107.14	121.92
1	B	443	SER	C-N-CA	-8.54	107.14	121.92

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	229	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	7396	0	7381	91	0
1	B	7354	0	7328	91	0
2	A	9	0	3	0	0
2	B	9	0	3	0	0
3	A	4	0	6	0	0
3	B	4	0	6	1	0
4	A	10	0	0	1	0
4	B	10	0	0	1	0
5	A	35	0	0	0	0
5	B	21	0	0	0	0
All	All	14852	0	14727	178	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 6.

The worst 5 of 178 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:806:LEU:HA	1:A:809:MET:HE2	1.31	1.12
1:A:240:MET:HE2	1:A:284:MET:HE1	1.46	0.98
1:A:748:MET:HE1	1:A:948:GLU:HB3	1.47	0.96
1:B:273:TYR:OH	1:B:411:GLU:OE2	1.90	0.90
1:A:806:LEU:HD23	1:A:809:MET:HE1	1.51	0.88

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	912/972 (94%)	879 (96%)	30 (3%)	3 (0%)	36	59
1	B	905/972 (93%)	873 (96%)	27 (3%)	5 (1%)	21	42
All	All	1817/1944 (94%)	1752 (96%)	57 (3%)	8 (0%)	30	52

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	752	SER
1	B	752	SER
1	A	22	GLY
1	B	22	GLY
1	B	498	ASP

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	802/848 (95%)	759 (95%)	43 (5%)	20	43
1	B	797/848 (94%)	756 (95%)	41 (5%)	21	46
All	All	1599/1696 (94%)	1515 (95%)	84 (5%)	21	44

5 of 84 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	225	ILE

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Mol	Chain	Res	Type
1	B	629	LEU
1	B	327	MET
1	B	529	SER
1	B	740	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 21 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	358	GLN
1	B	653	GLN
1	B	963	GLN
1	B	673	GLN
1	B	589	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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
4	SO4	B	1970	-	4,4,4	0.63	0	6,6,6	0.61	0
2	ASP	A	1967	-	7,8,8	1.43	2 (28%)	6,10,10	1.05	0
4	SO4	A	1969	-	4,4,4	0.39	0	6,6,6	1.26	1 (16%)
4	SO4	A	1970	-	4,4,4	0.55	0	6,6,6	0.32	0
3	EDO	B	1968	-	3,3,3	0.61	0	2,2,2	0.24	0
2	ASP	B	1967	-	7,8,8	1.32	0	6,10,10	1.22	1 (16%)
3	EDO	A	1968	-	3,3,3	0.89	0	2,2,2	0.44	0
4	SO4	B	1969	-	4,4,4	0.56	0	6,6,6	0.67	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
3	EDO	A	1968	-	-	1/1/1/1	-
2	ASP	A	1967	-	-	0/8/8/8	-
3	EDO	B	1968	-	-	1/1/1/1	-
2	ASP	B	1967	-	-	0/8/8/8	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1967	ASP	OXT-C	-2.23	1.23	1.30
2	A	1967	ASP	CB-CG	-2.05	1.45	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1967	ASP	OD1-CG-CB	-2.28	115.86	122.84
4	A	1969	SO4	O4-S-O3	2.00	119.59	108.54

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1968	EDO	O1-C1-C2-O2
3	B	1968	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1969	SO4	1	0
3	B	1968	EDO	1	0
4	B	1969	SO4	1	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	920/972 (94%)	-0.27	24 (2%) 57 54	23, 39, 78, 116	0
1	B	915/972 (94%)	-0.12	39 (4%) 40 36	29, 46, 88, 135	0
All	All	1835/1944 (94%)	-0.19	63 (3%) 48 44	23, 43, 83, 135	0

The worst 5 of 63 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	6	LEU	5.3
1	A	24	VAL	5.1
1	A	199	ASP	5.1
1	B	24	VAL	4.8
1	B	754	PRO	4.8

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	B	1970	5/5	0.87	0.18	60,72,86,87	0
4	SO4	A	1970	5/5	0.90	0.12	63,64,80,87	0
4	SO4	B	1969	5/5	0.90	0.15	55,64,69,76	0
3	EDO	A	1968	4/4	0.90	0.14	27,29,32,37	0
3	EDO	B	1968	4/4	0.94	0.12	31,36,37,39	0
4	SO4	A	1969	5/5	0.95	0.10	40,52,55,59	0
2	ASP	A	1967	9/9	0.95	0.09	37,39,46,47	0
2	ASP	B	1967	9/9	0.96	0.08	34,41,46,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.