



# wwPDB X-ray Structure Validation Summary Report ⓘ

Dec 10, 2022 – 06:37 PM EST

PDB ID : 1TD9  
Title : Crystal Structure of a Phosphotransacetylase from *Bacillus subtilis*  
Authors : Xu, Q.S.; Jancarik, J.; Yokota, H.; Kim, R.; Kim, S.-H.; Berkeley Structural Genomics Center (BSGC)  
Deposited on : 2004-05-21  
Resolution : 2.75 Å(reported)

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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>.

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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	:	2.31.2
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)	:	2.31.2

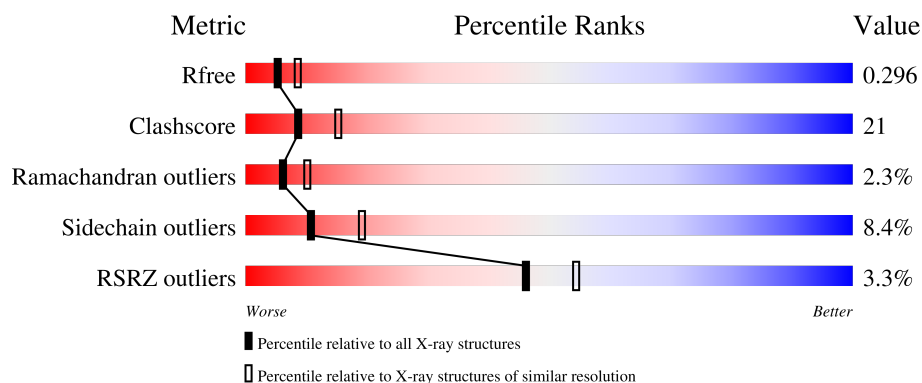
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.75 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	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)

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	329	<div> <div>0%</div> <div>63% 31% 5% .</div> </div>
1	B	329	<div> <div>2%</div> <div>60% 33% 6% .</div> </div>
1	C	329	<div> <div>2%</div> <div>61% 33% 5% .</div> </div>
1	D	329	<div> <div>2%</div> <div>57% 36% 6% .</div> </div>
1	E	329	<div> <div>2%</div> <div>61% 31% 6% .</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	329	<div><div></div><div>11%</div><div>59%</div><div>33%</div><div>6%</div><div></div></div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14916 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 Phosphate acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	0	0	0
			2445	1545	407	484	9			
1	B	324	Total	C	N	O	S	0	0	0
			2449	1547	408	485	9			
1	C	323	Total	C	N	O	S	0	0	0
			2445	1545	407	484	9			
1	D	325	Total	C	N	O	S	0	0	0
			2453	1549	409	486	9			
1	E	323	Total	C	N	O	S	0	0	0
			2445	1545	407	484	9			
1	F	323	Total	C	N	O	S	0	0	0
			2445	1545	407	484	9			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	cloning artifact	UNP P39646
A	-4	GLY	-	cloning artifact	UNP P39646
A	-3	GLY	-	cloning artifact	UNP P39646
A	-2	GLY	-	cloning artifact	UNP P39646
A	-1	GLY	-	cloning artifact	UNP P39646
A	0	GLY	-	cloning artifact	UNP P39646
A	1	MET	-	initiating methionine	UNP P39646
B	-5	GLY	-	cloning artifact	UNP P39646
B	-4	GLY	-	cloning artifact	UNP P39646
B	-3	GLY	-	cloning artifact	UNP P39646
B	-2	GLY	-	cloning artifact	UNP P39646
B	-1	GLY	-	cloning artifact	UNP P39646
B	0	GLY	-	cloning artifact	UNP P39646
B	1	MET	-	initiating methionine	UNP P39646
C	-5	GLY	-	cloning artifact	UNP P39646
C	-4	GLY	-	cloning artifact	UNP P39646
C	-3	GLY	-	cloning artifact	UNP P39646

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	GLY	-	cloning artifact	UNP P39646
C	-1	GLY	-	cloning artifact	UNP P39646
C	0	GLY	-	cloning artifact	UNP P39646
C	1	MET	-	initiating methionine	UNP P39646
D	-5	GLY	-	cloning artifact	UNP P39646
D	-4	GLY	-	cloning artifact	UNP P39646
D	-3	GLY	-	cloning artifact	UNP P39646
D	-2	GLY	-	cloning artifact	UNP P39646
D	-1	GLY	-	cloning artifact	UNP P39646
D	0	GLY	-	cloning artifact	UNP P39646
D	1	MET	-	initiating methionine	UNP P39646
E	-5	GLY	-	cloning artifact	UNP P39646
E	-4	GLY	-	cloning artifact	UNP P39646
E	-3	GLY	-	cloning artifact	UNP P39646
E	-2	GLY	-	cloning artifact	UNP P39646
E	-1	GLY	-	cloning artifact	UNP P39646
E	0	GLY	-	cloning artifact	UNP P39646
E	1	MET	-	initiating methionine	UNP P39646
F	-5	GLY	-	cloning artifact	UNP P39646
F	-4	GLY	-	cloning artifact	UNP P39646
F	-3	GLY	-	cloning artifact	UNP P39646
F	-2	GLY	-	cloning artifact	UNP P39646
F	-1	GLY	-	cloning artifact	UNP P39646
F	0	GLY	-	cloning artifact	UNP P39646
F	1	MET	-	initiating methionine	UNP P39646

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		

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

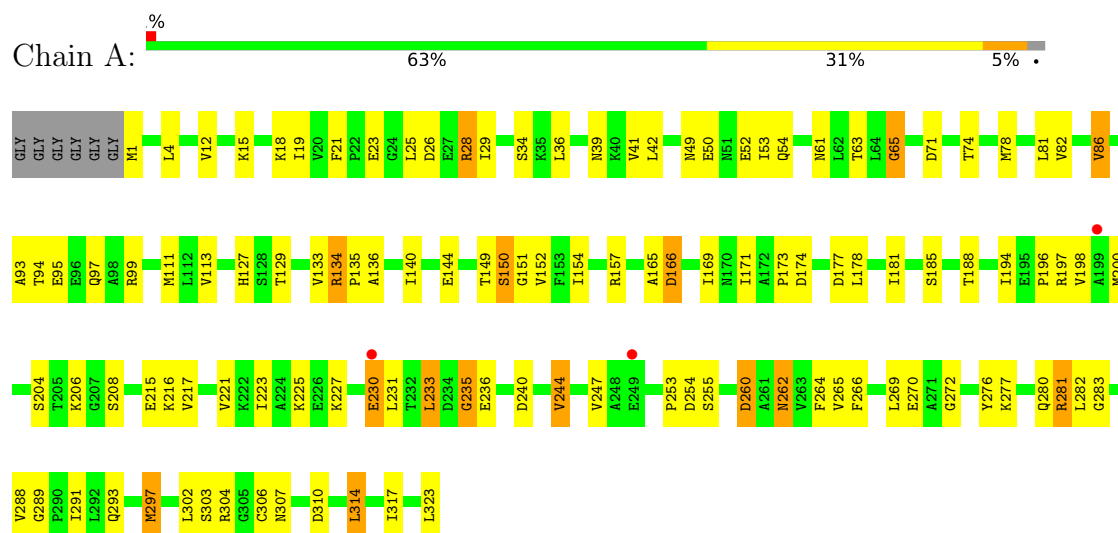
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	25	Total	O	0	0
			25	25		
3	B	18	Total	O	0	0
			18	18		
3	C	26	Total	O	0	0
			26	26		
3	D	18	Total	O	0	0
			18	18		
3	E	20	Total	O	0	0
			20	20		
3	F	7	Total	O	0	0
			7	7		

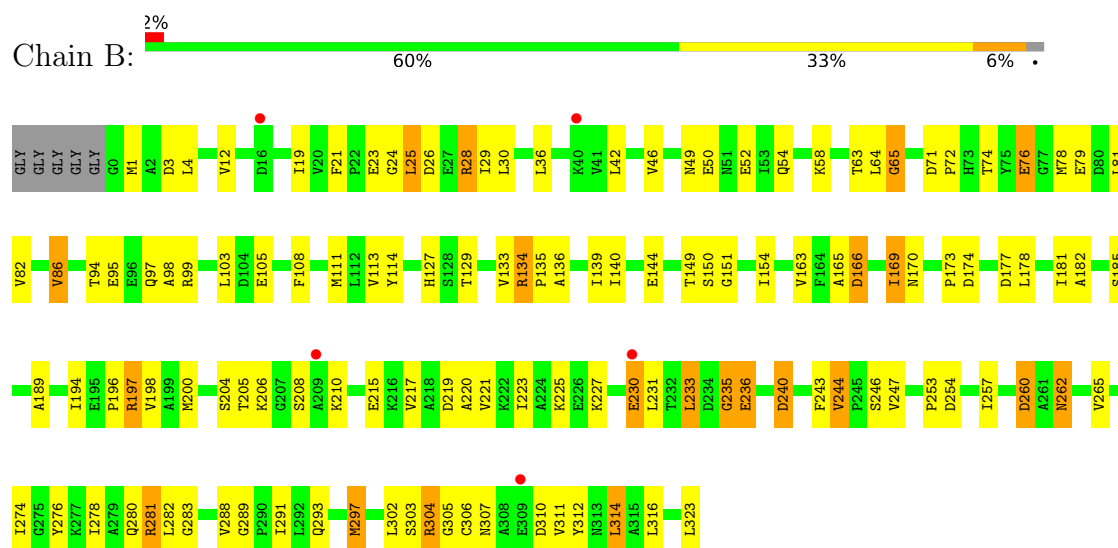
### 3 Residue-property plots [i](#)

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: Phosphate acetyltransferase



#### • Molecule 1: Phosphate acetyltransferase

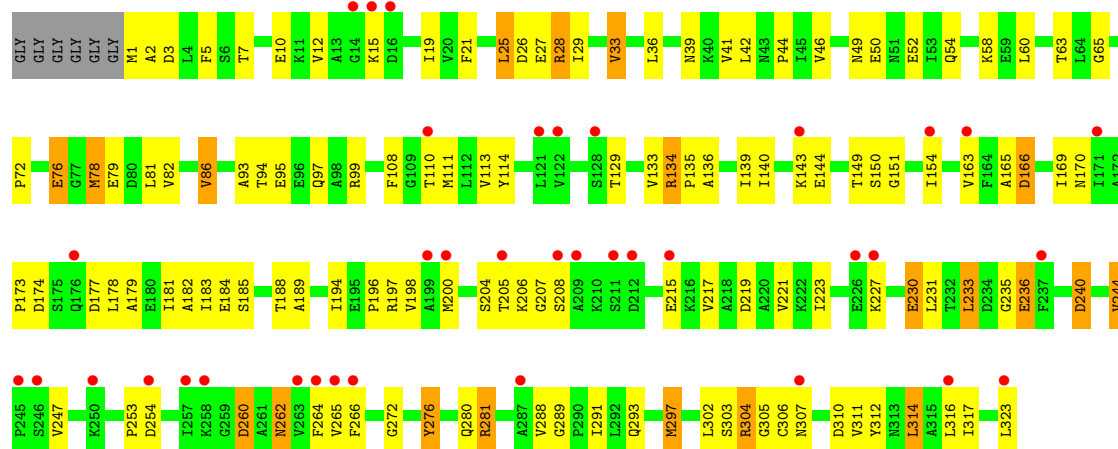


#### • Molecule 1: Phosphate acetyltransferase









## 4 Data and refinement statistics

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	185.21Å 185.21Å 259.41Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.99 – 2.75 19.99 – 2.75	Depositor EDS
% Data completeness (in resolution range)	94.5 (19.99-2.75) 94.6 (19.99-2.75)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	0.12	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.29 (at 2.75Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.268 , 0.297 0.267 , 0.296	Depositor DCC
$R_{free}$ test set	1958 reflections (3.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	45.6	Xtriage
Anisotropy	0.695	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 56.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	14916	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 53.77 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.0172e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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: 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	0.48	0/2479	0.75	2/3346 (0.1%)
1	B	0.45	0/2483	0.75	3/3351 (0.1%)
1	C	0.47	0/2479	0.74	1/3346 (0.0%)
1	D	0.44	0/2487	0.74	3/3356 (0.1%)
1	E	0.48	0/2479	0.74	2/3346 (0.1%)
1	F	0.45	0/2479	0.72	2/3346 (0.1%)
All	All	0.46	0/14886	0.74	13/20091 (0.1%)

There are no bond length outliers.

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	260	ASP	CB-CG-OD2	6.96	124.57	118.30
1	D	260	ASP	CB-CG-OD2	6.94	124.55	118.30
1	F	260	ASP	CB-CG-OD2	6.58	124.22	118.30
1	E	260	ASP	CB-CG-OD2	6.54	124.19	118.30
1	C	260	ASP	CB-CG-OD2	6.15	123.83	118.30

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	2445	0	2468	97	0
1	B	2449	0	2471	112	0
1	C	2445	0	2468	104	0
1	D	2453	0	2474	117	0
1	E	2445	0	2468	106	0
1	F	2445	0	2468	104	0
2	A	25	0	0	0	0
2	B	20	0	0	0	0
2	C	25	0	0	0	0
2	D	20	0	0	1	0
2	E	20	0	0	0	0
2	F	10	0	0	0	0
3	A	25	0	0	2	0
3	B	18	0	0	4	0
3	C	26	0	0	3	0
3	D	18	0	0	2	0
3	E	20	0	0	0	0
3	F	7	0	0	1	0
All	All	14916	0	14817	610	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 21.

The worst 5 of 610 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:165:ALA:HB3	1:A:265:VAL:HA	1.30	1.14
1:C:165:ALA:HB3	1:C:265:VAL:HA	1.30	1.13
1:D:165:ALA:HB3	1:D:265:VAL:HA	1.32	1.11
1:B:165:ALA:HB3	1:B:265:VAL:HA	1.34	1.08
1:E:165:ALA:HB3	1:E:265:VAL:HA	1.33	1.07

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	321/329 (98%)	289 (90%)	26 (8%)	6 (2%)	8	14
1	B	322/329 (98%)	288 (89%)	26 (8%)	8 (2%)	5	9
1	C	321/329 (98%)	290 (90%)	25 (8%)	6 (2%)	8	14
1	D	323/329 (98%)	287 (89%)	28 (9%)	8 (2%)	5	9
1	E	321/329 (98%)	285 (89%)	28 (9%)	8 (2%)	5	9
1	F	321/329 (98%)	281 (88%)	31 (10%)	9 (3%)	5	7
All	All	1929/1974 (98%)	1720 (89%)	164 (8%)	45 (2%)	6	10

5 of 45 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	65	GLY
1	B	144	GLU
1	D	144	GLU
1	F	2	ALA
1	F	144	GLU

### 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	257/257 (100%)	237 (92%)	20 (8%)	12	22
1	B	257/257 (100%)	235 (91%)	22 (9%)	10	18
1	C	257/257 (100%)	236 (92%)	21 (8%)	11	20
1	D	257/257 (100%)	236 (92%)	21 (8%)	11	20
1	E	257/257 (100%)	235 (91%)	22 (9%)	10	18
1	F	257/257 (100%)	234 (91%)	23 (9%)	9	17
All	All	1542/1542 (100%)	1413 (92%)	129 (8%)	11	19

5 of 129 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	149	THR
1	F	227	LYS
1	C	134	ARG
1	C	86	VAL
1	F	233	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 35 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	293	GLN
1	F	49	ASN
1	F	176	GLN
1	C	127	HIS
1	C	49	ASN

### 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](#)

24 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	SO4	B	407	-	4,4,4	2.11	1 (25%)	6,6,6	1.01	1 (16%)
2	SO4	D	416	-	4,4,4	2.24	1 (25%)	6,6,6	1.30	2 (33%)
2	SO4	F	423	-	4,4,4	2.33	1 (25%)	6,6,6	0.70	0
2	SO4	A	401	-	4,4,4	2.47	1 (25%)	6,6,6	0.97	1 (16%)
2	SO4	C	412	-	4,4,4	2.44	1 (25%)	6,6,6	0.75	0
2	SO4	E	421	-	4,4,4	2.40	1 (25%)	6,6,6	0.82	0
2	SO4	E	419	-	4,4,4	2.55	1 (25%)	6,6,6	1.05	1 (16%)
2	SO4	D	413	-	4,4,4	2.03	1 (25%)	6,6,6	0.69	0
2	SO4	C	411	-	4,4,4	2.02	1 (25%)	6,6,6	1.18	1 (16%)
2	SO4	A	403	-	4,4,4	2.52	1 (25%)	6,6,6	0.77	0
2	SO4	C	417	-	4,4,4	2.35	1 (25%)	6,6,6	0.85	0
2	SO4	A	405	-	4,4,4	2.17	1 (25%)	6,6,6	0.87	0
2	SO4	E	420	-	4,4,4	2.19	1 (25%)	6,6,6	1.33	2 (33%)
2	SO4	C	410	-	4,4,4	2.39	1 (25%)	6,6,6	0.89	0
2	SO4	D	415	-	4,4,4	2.65	1 (25%)	6,6,6	0.74	0
2	SO4	B	406	-	4,4,4	2.31	1 (25%)	6,6,6	0.66	0
2	SO4	D	418	-	4,4,4	2.66	1 (25%)	6,6,6	0.70	0
2	SO4	A	402	-	4,4,4	1.90	1 (25%)	6,6,6	1.27	2 (33%)
2	SO4	B	404	-	4,4,4	2.41	1 (25%)	6,6,6	0.79	0
2	SO4	B	409	-	4,4,4	2.51	1 (25%)	6,6,6	0.86	0
2	SO4	A	408	-	4,4,4	2.49	2 (50%)	6,6,6	1.21	1 (16%)
2	SO4	C	414	-	4,4,4	2.53	1 (25%)	6,6,6	0.79	0
2	SO4	F	424	-	4,4,4	2.09	1 (25%)	6,6,6	1.08	0
2	SO4	E	422	-	4,4,4	2.16	1 (25%)	6,6,6	1.01	1 (16%)

The worst 5 of 25 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	418	SO4	O1-S	5.07	1.73	1.46
2	D	415	SO4	O1-S	4.87	1.72	1.46
2	E	419	SO4	O1-S	4.84	1.72	1.46
2	C	414	SO4	O1-S	4.80	1.72	1.46
2	A	403	SO4	O1-S	4.80	1.72	1.46

The worst 5 of 12 bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	420	SO4	O4-S-O3	2.56	119.99	109.06
2	D	416	SO4	O4-S-O3	2.34	119.04	109.06
2	A	408	SO4	O4-S-O3	2.33	118.99	109.06
2	C	411	SO4	O4-S-O3	2.22	118.55	109.06
2	A	402	SO4	O4-S-O3	2.16	118.30	109.06

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	418	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, 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	323/329 (98%)	-0.23	3 (0%) 84 89	21, 42, 70, 90	0
1	B	324/329 (98%)	-0.05	5 (1%) 73 81	22, 48, 78, 95	0
1	C	323/329 (98%)	-0.13	5 (1%) 73 81	18, 43, 78, 99	0
1	D	325/329 (98%)	0.03	8 (2%) 57 66	18, 49, 81, 96	0
1	E	323/329 (98%)	0.02	6 (1%) 66 75	20, 49, 76, 96	0
1	F	323/329 (98%)	0.76	37 (11%) 4 5	34, 75, 99, 106	0
All	All	1941/1974 (98%)	0.07	64 (3%) 46 54	18, 50, 87, 106	0

The worst 5 of 64 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	254	ASP	3.8
1	F	14	GLY	3.4
1	F	199	ALA	3.3
1	C	2	ALA	3.3
1	F	287	ALA	3.2

### 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 ⓘ

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(Å <sup>2</sup> )	Q<0.9
2	SO4	D	416	5/5	0.84	0.32	66,70,79,83	0
2	SO4	C	412	5/5	0.86	0.41	69,77,83,89	0
2	SO4	E	420	5/5	0.88	0.34	67,70,75,80	0
2	SO4	E	421	5/5	0.88	0.29	65,66,76,84	0
2	SO4	C	414	5/5	0.89	0.22	72,72,85,88	0
2	SO4	E	419	5/5	0.89	0.35	60,71,79,83	0
2	SO4	A	403	5/5	0.90	0.30	68,71,78,84	0
2	SO4	B	407	5/5	0.90	0.36	66,76,83,85	0
2	SO4	B	409	5/5	0.90	0.26	61,66,83,84	0
2	SO4	A	408	5/5	0.91	0.29	70,71,80,82	0
2	SO4	D	415	5/5	0.91	0.31	67,69,84,84	0
2	SO4	E	422	5/5	0.91	0.29	55,69,74,75	0
2	SO4	F	424	5/5	0.91	0.32	68,71,83,83	0
2	SO4	D	418	5/5	0.92	0.23	58,62,75,77	0
2	SO4	B	406	5/5	0.92	0.31	76,76,86,88	0
2	SO4	F	423	5/5	0.92	0.33	74,78,84,86	0
2	SO4	D	413	5/5	0.92	0.23	72,74,80,83	0
2	SO4	C	410	5/5	0.93	0.30	62,73,75,76	0
2	SO4	A	401	5/5	0.94	0.29	57,76,76,77	0
2	SO4	C	411	5/5	0.94	0.25	64,72,76,77	0
2	SO4	A	402	5/5	0.94	0.27	59,63,66,69	0
2	SO4	B	404	5/5	0.94	0.25	70,76,77,79	0
2	SO4	C	417	5/5	0.94	0.22	66,78,80,86	0
2	SO4	A	405	5/5	0.95	0.21	63,67,75,77	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.