



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 10:43 AM UTC

PDB ID : 2FEN / pdb\_00002fen  
Title : 3-carboxy-cis,cis-muconate lactonizing enzyme from Agrobacterium radiobacter S2  
Authors : Lehtio, L.; Goldman, A.  
Deposited on : 2005-12-16  
Resolution : 2.60 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](mailto: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>.

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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

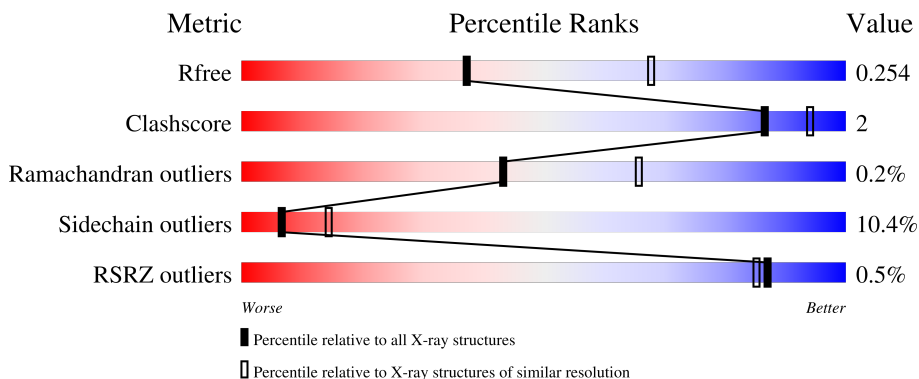
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	359	
1	B	359	
1	C	359	
1	D	359	
1	E	359	

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Mol	Chain	Length	Quality of chain
1	F	359	 82% 10% • 6%
1	G	359	 83% 9% • 6%
1	H	359	 83% 9% • 6%
1	I	359	 82% 9% • 6%
1	J	359	 80% 12% • 6%
1	K	359	 82% 11% • 6%
1	L	359	 83% 9% • 6%

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 31433 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 3-carboxy-cis,cis-muconate lactonizing enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	342	2574	1615	457	493	9	0	2	0
1	B	339	2545	1598	452	486	9	0	1	0
1	D	343	2564	1610	456	489	9	0	0	0
1	C	338	2543	1597	450	487	9	0	1	0
1	E	337	2544	1598	450	487	9	0	2	0
1	F	339	2537	1594	450	484	9	0	0	0
1	H	337	2552	1602	452	489	9	0	3	0
1	G	337	2551	1602	451	489	9	0	3	0
1	I	337	2544	1598	450	487	9	0	2	0
1	J	339	2545	1598	452	486	9	0	1	0
1	L	337	2552	1602	452	489	9	0	3	0
1	K	336	2532	1592	448	483	9	0	1	0

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

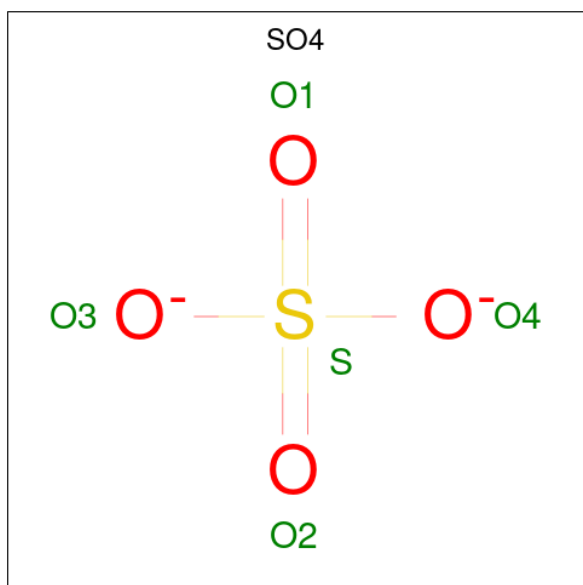
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Cl	0	0
			1	1		
2	C	2	Total	Cl	0	0
			2	2		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	I	1	Total Cl 1 1	0	0
2	J	1	Total Cl 1 1	0	0

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	H	1	Total O S 5 4 1	0	0
3	G	1	Total O S 5 4 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	79	Total O 79 79	0	0
4	B	79	Total O 79 79	0	0
4	D	98	Total O 98 98	0	0

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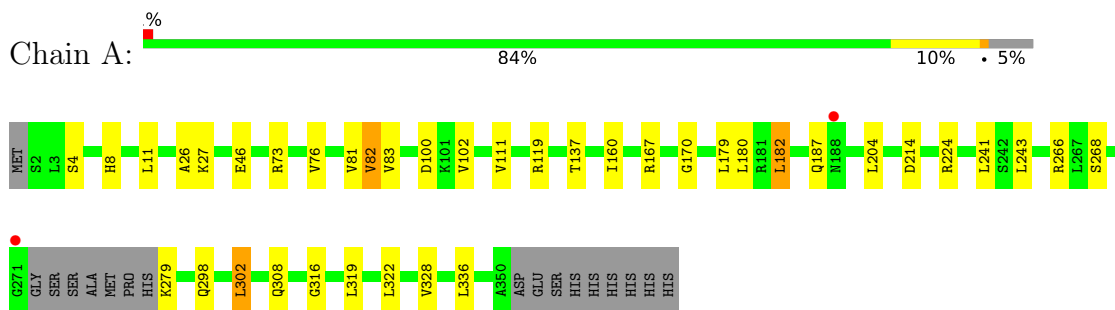
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
4	C	103	Total 103	O 103	0	0
4	E	45	Total 45	O 45	0	0
4	F	72	Total 72	O 72	0	0
4	H	74	Total 74	O 74	0	0
4	G	72	Total 72	O 72	0	0
4	I	50	Total 50	O 50	0	0
4	J	53	Total 53	O 53	0	0
4	L	30	Total 30	O 30	0	0
4	K	70	Total 70	O 70	0	0

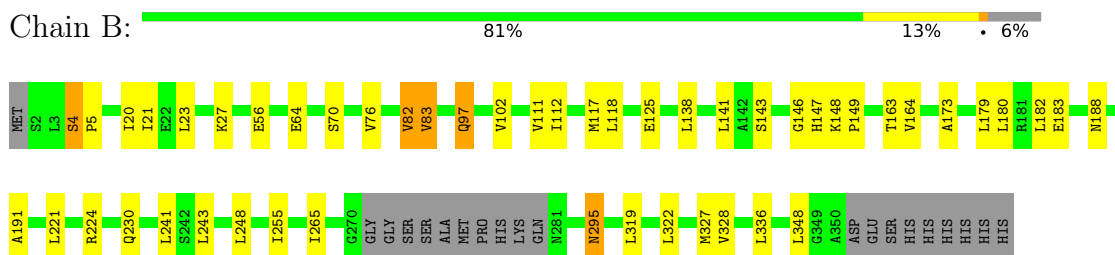
### 3 Residue-property plots [i](#)

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.

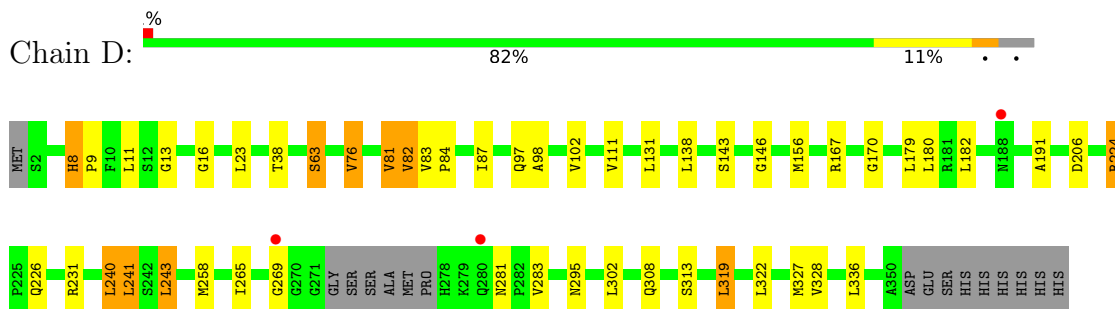
- Molecule 1: 3-carboxy-cis,cis-muconate lactonizing enzyme



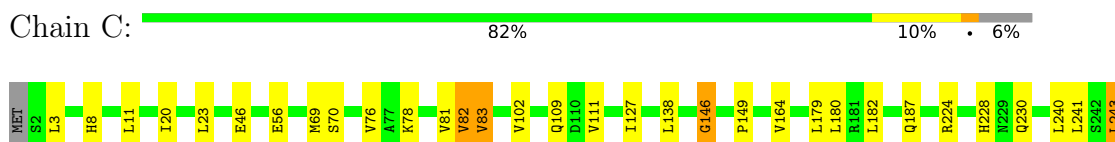
- Molecule 1: 3-carboxy-cis,cis-muconate lactonizing enzyme

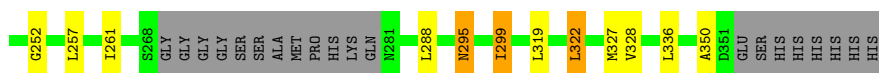


- Molecule 1: 3-carboxy-cis,cis-muconate lactonizing enzyme



- Molecule 1: 3-carboxy-cis,cis-muconate lactonizing enzyme





- Molecule 1: 3-carboxy-cis,cis-muconate lactonizing enzyme

Chain E: 82% 11% • 6%



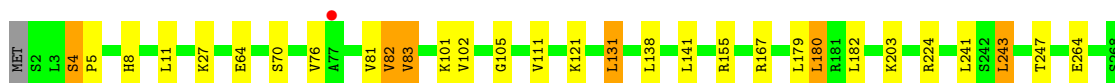
- Molecule 1: 3-carboxy-cis,cis-muconate lactonizing enzyme

Chain F: 82% 10% • 6%



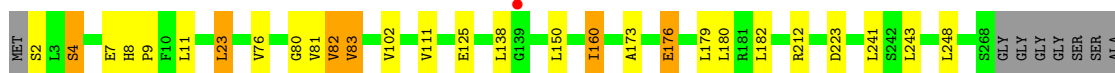
- Molecule 1: 3-carboxy-cis,cis-muconate lactonizing enzyme

Chain H: 83% 9% • 6%



- Molecule 1: 3-carboxy-cis,cis-muconate lactonizing enzyme

Chain G: 83% 9% • 6%



- Molecule 1: 3-carboxy-cis,cis-muconate lactonizing enzyme

Chain I: 82% 9% • 6%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.03Å 205.32Å 235.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 2.60 19.99 – 2.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.99-2.60) 94.6 (19.99-2.60)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.23 (at 2.59Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.200 , 0.262 (Not available) , 0.254	Depositor DCC
$R_{free}$ test set	6657 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.6	Xtrriage
Anisotropy	0.076	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 46.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	31433	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.42% of the height of the origin peak. No significant pseudotranslation is detected.*

<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, CL

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	0.69	0/2611	1.11	10/3531 (0.3%)
1	B	0.70	0/2582	1.10	8/3494 (0.2%)
1	C	0.76	1/2580 (0.0%)	1.12	4/3492 (0.1%)
1	D	0.73	0/2601	1.13	9/3518 (0.3%)
1	E	0.58	0/2581	1.03	8/3493 (0.2%)
1	F	0.63	0/2574	1.07	13/3483 (0.4%)
1	G	0.64	0/2589	1.04	7/3504 (0.2%)
1	H	0.63	0/2589	1.01	4/3504 (0.1%)
1	I	0.69	1/2581 (0.0%)	1.07	9/3493 (0.3%)
1	J	0.89	11/2582 (0.4%)	1.06	2/3494 (0.1%)
1	K	0.62	1/2569 (0.0%)	1.05	7/3477 (0.2%)
1	L	0.66	1/2589 (0.0%)	1.04	7/3504 (0.2%)
All	All	0.69	15/31028 (0.0%)	1.07	88/41987 (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	G	0	1

The worst 5 of 15 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	52	ASP	CG-OD1	13.48	1.50	1.25
1	J	52	ASP	CG-OD2	9.33	1.43	1.25
1	J	64	GLU	CD-OE2	7.68	1.40	1.25
1	J	53	ASP	CG-OD1	7.44	1.39	1.25
1	J	56	GLU	CD-OE2	7.30	1.39	1.25

The worst 5 of 88 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	147	HIS	N-CA-C	-9.67	102.57	114.75
1	L	82	VAL	CB-CA-C	-9.42	99.91	111.97
1	I	82	VAL	CB-CA-C	-9.33	99.74	112.24
1	E	4	SER	CA-C-N	8.35	129.16	119.47
1	E	4	SER	C-N-CA	8.35	129.16	119.47

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	G	281	ASN	Peptide

## 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	2574	0	2575	10	0
1	B	2545	0	2547	15	0
1	C	2543	0	2544	18	0
1	D	2564	0	2568	16	0
1	E	2544	0	2545	8	0
1	F	2537	0	2542	13	0
1	G	2551	0	2550	11	0
1	H	2552	0	2550	12	0
1	I	2544	0	2545	14	0
1	J	2545	0	2547	9	0
1	K	2532	0	2536	8	0
1	L	2552	0	2550	10	0
2	A	1	0	0	0	0
2	C	2	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	0	0
3	G	5	0	0	0	0
3	H	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	79	0	0	0	0
4	B	79	0	0	0	0
4	C	103	0	0	0	0
4	D	98	0	0	0	0
4	E	45	0	0	0	0
4	F	72	0	0	1	0
4	G	72	0	0	0	0
4	H	74	0	0	0	0
4	I	50	0	0	0	0
4	J	53	0	0	0	0
4	K	70	0	0	1	0
4	L	30	0	0	0	0
All	All	31433	0	30599	128	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 2.

The worst 5 of 128 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:C:295:ASN:HD21	1:C:327:MET:HA	1.40	0.87
1:D:170:GLY:HA2	1:C:230:GLN:HE21	1.50	0.76
1:A:170:GLY:HA2	1:B:230:GLN:HE21	1.58	0.67
1:F:283:VAL:HG11	1:G:80:GLY:HA3	1.77	0.65
1:J:58:ILE:HD11	1:J:101:LYS:HB2	1.79	0.65

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	340/359 (95%)	328 (96%)	12 (4%)	0	100	100
1	B	336/359 (94%)	326 (97%)	10 (3%)	0	100	100
1	C	335/359 (93%)	318 (95%)	17 (5%)	0	100	100
1	D	339/359 (94%)	326 (96%)	11 (3%)	2 (1%)	21	42
1	E	335/359 (93%)	321 (96%)	12 (4%)	2 (1%)	21	42
1	F	335/359 (93%)	326 (97%)	9 (3%)	0	100	100
1	G	336/359 (94%)	322 (96%)	14 (4%)	0	100	100
1	H	336/359 (94%)	321 (96%)	15 (4%)	0	100	100
1	I	335/359 (93%)	322 (96%)	12 (4%)	1 (0%)	36	58
1	J	336/359 (94%)	315 (94%)	20 (6%)	1 (0%)	36	58
1	K	333/359 (93%)	321 (96%)	11 (3%)	1 (0%)	36	58
1	L	336/359 (94%)	323 (96%)	13 (4%)	0	100	100
All	All	4032/4308 (94%)	3869 (96%)	156 (4%)	7 (0%)	43	66

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	63	SER
1	E	189	GLY
1	J	99	ALA
1	I	17	ASP
1	K	64	GLU

### 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	258/271 (95%)	235 (91%)	23 (9%)	9	20
1	B	255/271 (94%)	227 (89%)	28 (11%)	6	13
1	C	256/271 (94%)	228 (89%)	28 (11%)	6	13
1	D	256/271 (94%)	229 (90%)	27 (10%)	6	14

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	256/271 (94%)	228 (89%)	28 (11%)	6	13
1	F	254/271 (94%)	228 (90%)	26 (10%)	7	15
1	G	257/271 (95%)	232 (90%)	25 (10%)	8	17
1	H	257/271 (95%)	232 (90%)	25 (10%)	8	17
1	I	256/271 (94%)	233 (91%)	23 (9%)	9	20
1	J	255/271 (94%)	221 (87%)	34 (13%)	4	8
1	K	254/271 (94%)	227 (89%)	27 (11%)	6	14
1	L	257/271 (95%)	233 (91%)	24 (9%)	8	18
All	All	3071/3252 (94%)	2753 (90%)	318 (10%)	7	14

5 of 318 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	328	VAL
1	L	295	ASN
1	J	64	GLU
1	J	243	LEU
1	K	102	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 58 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	207	ASN
1	K	284	ASN
1	G	187	GLN
1	K	281	ASN
1	L	157	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](#)

Of 9 ligands modelled in this entry, 5 are 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
3	SO4	B	1001	-	4,4,4	0.26	0	6,6,6	0.22	0
3	SO4	G	1004	-	4,4,4	0.26	0	6,6,6	0.18	0
3	SO4	H	1002	-	4,4,4	0.29	0	6,6,6	0.14	0
3	SO4	C	1003	-	4,4,4	0.21	0	6,6,6	0.10	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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 [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	342/359 (95%)	-0.39	2 (0%) 85 83	14, 28, 36, 52	3 (0%)
1	B	339/359 (94%)	-0.45	0 100 100	17, 28, 36, 52	1 (0%)
1	C	338/359 (94%)	-0.40	0 100 100	17, 27, 34, 39	1 (0%)
1	D	343/359 (95%)	-0.41	3 (0%) 81 78	17, 28, 37, 50	1 (0%)
1	E	337/359 (93%)	-0.20	1 (0%) 90 88	15, 28, 34, 37	3 (0%)
1	F	339/359 (94%)	-0.41	0 100 100	21, 28, 36, 45	0
1	G	337/359 (93%)	-0.22	1 (0%) 90 88	13, 28, 35, 41	3 (0%)
1	H	337/359 (93%)	-0.16	1 (0%) 90 88	14, 28, 35, 39	3 (0%)
1	I	337/359 (93%)	0.04	4 (1%) 76 73	14, 28, 34, 38	3 (0%)
1	J	339/359 (94%)	0.12	5 (1%) 72 68	15, 29, 36, 41	1 (0%)
1	K	336/359 (93%)	-0.06	1 (0%) 90 88	16, 29, 36, 40	1 (0%)
1	L	337/359 (93%)	-0.02	2 (0%) 85 83	16, 28, 36, 40	3 (0%)
All	All	4061/4308 (94%)	-0.21	20 (0%) 87 85	13, 28, 36, 52	23 (0%)

The worst 5 of 20 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	188	ASN	4.8
1	I	161	GLY	4.2
1	A	188	ASN	4.1
1	I	188	ASN	3.5
1	D	188	ASN	3.4

### 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, 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( $\text{\AA}^2$ )	Q<0.9
3	SO4	H	1002	5/5	0.80	0.14	87,87,87,87	0
3	SO4	G	1004	5/5	0.89	0.18	65,65,66,67	0
3	SO4	B	1001	5/5	0.92	0.13	54,55,56,56	0
2	CL	C	2004	1/1	0.93	0.08	40,40,40,40	0
3	SO4	C	1003	5/5	0.94	0.15	64,65,65,65	0
2	CL	A	2002	1/1	0.95	0.09	31,31,31,31	0
2	CL	J	2001	1/1	0.95	0.09	35,35,35,35	0
2	CL	I	2005	1/1	0.99	0.09	24,24,24,24	0
2	CL	C	2003	1/1	0.99	0.06	25,25,25,25	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.