



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 29, 2017 – 12:48 PM EDT

PDB ID : 3KL2
Title : Crystal structure of a putative isochorismatase from *Streptomyces avermitilis*
Authors : Bonanno, J.B.; Dickey, M.; Bain, K.T.; Chang, S.; Ozyurt, S.; Wasserman, S.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : unknown
Resolution : 2.30 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030345

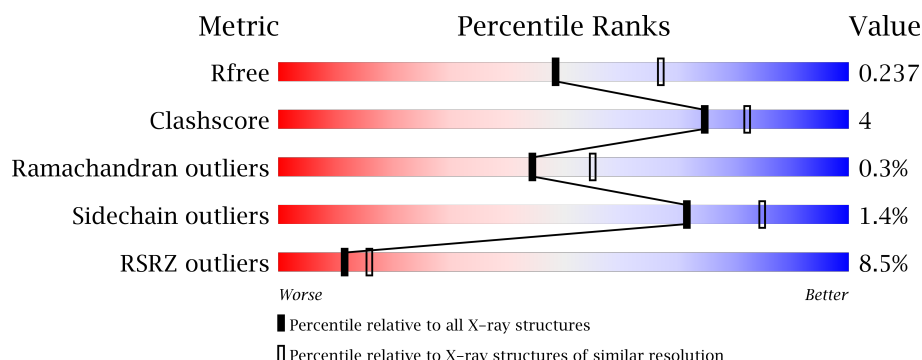
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	226	<div> <div>7%</div> <div> <div></div> <div>77%</div> <div>12%</div> <div>10%</div> </div> </div>
1	B	226	<div> <div>7%</div> <div> <div></div> <div>81%</div> <div>7%</div> <div>12%</div> </div> </div>
1	C	226	<div> <div>6%</div> <div> <div></div> <div>77%</div> <div>11%</div> <div>12%</div> </div> </div>
1	D	226	<div> <div>8%</div> <div> <div></div> <div>81%</div> <div>7%</div> <div>12%</div> </div> </div>
1	E	226	<div> <div>3%</div> <div> <div></div> <div>83%</div> <div>7%</div> <div>10%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	226	<div><div></div><div>12%</div><div>82%</div><div>7%</div><div>11%</div></div>
1	G	226	<div><div></div><div>3%</div><div>77%</div><div>10%</div><div>12%</div></div>
1	H	226	<div><div></div><div>5%</div><div>83%</div><div>6%</div><div>12%</div></div>
1	I	226	<div><div></div><div>15%</div><div>82%</div><div>6%</div><div>12%</div></div>
1	J	226	<div><div></div><div>16%</div><div>80%</div><div>8%</div><div>12%</div></div>
1	K	226	<div><div></div><div>5%</div><div>79%</div><div>9%</div><div>12%</div></div>
1	L	226	<div><div></div><div>5%</div><div>77%</div><div>12%</div><div>12%</div></div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 17957 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 Putative isochorismatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	S	0	0	0
			1491	945	251	286	9			
1	B	200	Total	C	N	O	S	0	0	0
			1460	926	242	283	9			
1	C	200	Total	C	N	O	S	0	0	0
			1475	935	246	285	9			
1	D	200	Total	C	N	O	S	0	0	0
			1488	941	250	288	9			
1	E	204	Total	C	N	O	S	0	0	0
			1513	956	258	290	9			
1	F	201	Total	C	N	O	S	0	0	0
			1493	946	251	287	9			
1	G	200	Total	C	N	O	S	0	0	0
			1488	941	250	288	9			
1	H	200	Total	C	N	O	S	0	0	0
			1485	940	250	286	9			
1	I	199	Total	C	N	O	S	0	0	0
			1456	922	247	278	9			
1	J	200	Total	C	N	O	S	0	0	0
			1469	931	246	283	9			
1	K	200	Total	C	N	O	S	0	0	0
			1475	935	247	284	9			
1	L	200	Total	C	N	O	S	0	0	0
			1480	937	249	285	9			

There are 132 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	18	MET	-	expression tag	UNP Q82NB5
A	19	SER	-	expression tag	UNP Q82NB5
A	20	LEU	-	expression tag	UNP Q82NB5
A	236	GLU	-	expression tag	UNP Q82NB5
A	237	GLY	-	expression tag	UNP Q82NB5

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Chain	Residue	Modelled	Actual	Comment	Reference
A	238	HIS	-	expression tag	UNP Q82NB5
A	239	HIS	-	expression tag	UNP Q82NB5
A	240	HIS	-	expression tag	UNP Q82NB5
A	241	HIS	-	expression tag	UNP Q82NB5
A	242	HIS	-	expression tag	UNP Q82NB5
A	243	HIS	-	expression tag	UNP Q82NB5
B	18	MET	-	expression tag	UNP Q82NB5
B	19	SER	-	expression tag	UNP Q82NB5
B	20	LEU	-	expression tag	UNP Q82NB5
B	236	GLU	-	expression tag	UNP Q82NB5
B	237	GLY	-	expression tag	UNP Q82NB5
B	238	HIS	-	expression tag	UNP Q82NB5
B	239	HIS	-	expression tag	UNP Q82NB5
B	240	HIS	-	expression tag	UNP Q82NB5
B	241	HIS	-	expression tag	UNP Q82NB5
B	242	HIS	-	expression tag	UNP Q82NB5
B	243	HIS	-	expression tag	UNP Q82NB5
C	18	MET	-	expression tag	UNP Q82NB5
C	19	SER	-	expression tag	UNP Q82NB5
C	20	LEU	-	expression tag	UNP Q82NB5
C	236	GLU	-	expression tag	UNP Q82NB5
C	237	GLY	-	expression tag	UNP Q82NB5
C	238	HIS	-	expression tag	UNP Q82NB5
C	239	HIS	-	expression tag	UNP Q82NB5
C	240	HIS	-	expression tag	UNP Q82NB5
C	241	HIS	-	expression tag	UNP Q82NB5
C	242	HIS	-	expression tag	UNP Q82NB5
C	243	HIS	-	expression tag	UNP Q82NB5
D	18	MET	-	expression tag	UNP Q82NB5
D	19	SER	-	expression tag	UNP Q82NB5
D	20	LEU	-	expression tag	UNP Q82NB5
D	236	GLU	-	expression tag	UNP Q82NB5
D	237	GLY	-	expression tag	UNP Q82NB5
D	238	HIS	-	expression tag	UNP Q82NB5
D	239	HIS	-	expression tag	UNP Q82NB5
D	240	HIS	-	expression tag	UNP Q82NB5
D	241	HIS	-	expression tag	UNP Q82NB5
D	242	HIS	-	expression tag	UNP Q82NB5
D	243	HIS	-	expression tag	UNP Q82NB5
E	18	MET	-	expression tag	UNP Q82NB5
E	19	SER	-	expression tag	UNP Q82NB5
E	20	LEU	-	expression tag	UNP Q82NB5

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Chain	Residue	Modelled	Actual	Comment	Reference
E	236	GLU	-	expression tag	UNP Q82NB5
E	237	GLY	-	expression tag	UNP Q82NB5
E	238	HIS	-	expression tag	UNP Q82NB5
E	239	HIS	-	expression tag	UNP Q82NB5
E	240	HIS	-	expression tag	UNP Q82NB5
E	241	HIS	-	expression tag	UNP Q82NB5
E	242	HIS	-	expression tag	UNP Q82NB5
E	243	HIS	-	expression tag	UNP Q82NB5
F	18	MET	-	expression tag	UNP Q82NB5
F	19	SER	-	expression tag	UNP Q82NB5
F	20	LEU	-	expression tag	UNP Q82NB5
F	236	GLU	-	expression tag	UNP Q82NB5
F	237	GLY	-	expression tag	UNP Q82NB5
F	238	HIS	-	expression tag	UNP Q82NB5
F	239	HIS	-	expression tag	UNP Q82NB5
F	240	HIS	-	expression tag	UNP Q82NB5
F	241	HIS	-	expression tag	UNP Q82NB5
F	242	HIS	-	expression tag	UNP Q82NB5
F	243	HIS	-	expression tag	UNP Q82NB5
G	18	MET	-	expression tag	UNP Q82NB5
G	19	SER	-	expression tag	UNP Q82NB5
G	20	LEU	-	expression tag	UNP Q82NB5
G	236	GLU	-	expression tag	UNP Q82NB5
G	237	GLY	-	expression tag	UNP Q82NB5
G	238	HIS	-	expression tag	UNP Q82NB5
G	239	HIS	-	expression tag	UNP Q82NB5
G	240	HIS	-	expression tag	UNP Q82NB5
G	241	HIS	-	expression tag	UNP Q82NB5
G	242	HIS	-	expression tag	UNP Q82NB5
G	243	HIS	-	expression tag	UNP Q82NB5
H	18	MET	-	expression tag	UNP Q82NB5
H	19	SER	-	expression tag	UNP Q82NB5
H	20	LEU	-	expression tag	UNP Q82NB5
H	236	GLU	-	expression tag	UNP Q82NB5
H	237	GLY	-	expression tag	UNP Q82NB5
H	238	HIS	-	expression tag	UNP Q82NB5
H	239	HIS	-	expression tag	UNP Q82NB5
H	240	HIS	-	expression tag	UNP Q82NB5
H	241	HIS	-	expression tag	UNP Q82NB5
H	242	HIS	-	expression tag	UNP Q82NB5
H	243	HIS	-	expression tag	UNP Q82NB5
I	18	MET	-	expression tag	UNP Q82NB5

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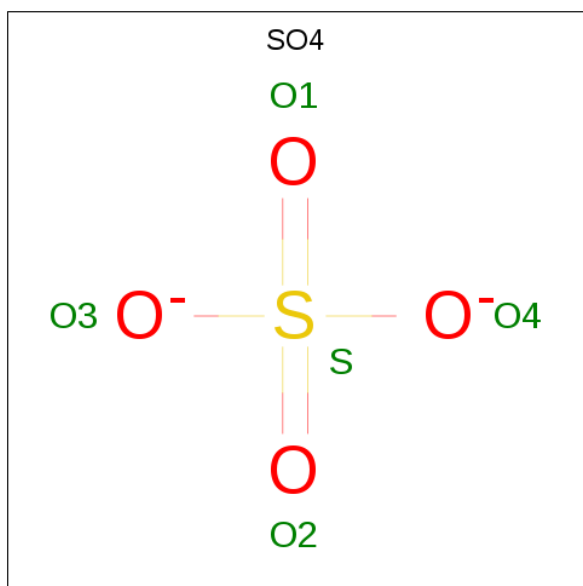
Chain	Residue	Modelled	Actual	Comment	Reference
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I	20	LEU	-	expression tag	UNP Q82NB5
I	236	GLU	-	expression tag	UNP Q82NB5
I	237	GLY	-	expression tag	UNP Q82NB5
I	238	HIS	-	expression tag	UNP Q82NB5
I	239	HIS	-	expression tag	UNP Q82NB5
I	240	HIS	-	expression tag	UNP Q82NB5
I	241	HIS	-	expression tag	UNP Q82NB5
I	242	HIS	-	expression tag	UNP Q82NB5
I	243	HIS	-	expression tag	UNP Q82NB5
J	18	MET	-	expression tag	UNP Q82NB5
J	19	SER	-	expression tag	UNP Q82NB5
J	20	LEU	-	expression tag	UNP Q82NB5
J	236	GLU	-	expression tag	UNP Q82NB5
J	237	GLY	-	expression tag	UNP Q82NB5
J	238	HIS	-	expression tag	UNP Q82NB5
J	239	HIS	-	expression tag	UNP Q82NB5
J	240	HIS	-	expression tag	UNP Q82NB5
J	241	HIS	-	expression tag	UNP Q82NB5
J	242	HIS	-	expression tag	UNP Q82NB5
J	243	HIS	-	expression tag	UNP Q82NB5
K	18	MET	-	expression tag	UNP Q82NB5
K	19	SER	-	expression tag	UNP Q82NB5
K	20	LEU	-	expression tag	UNP Q82NB5
K	236	GLU	-	expression tag	UNP Q82NB5
K	237	GLY	-	expression tag	UNP Q82NB5
K	238	HIS	-	expression tag	UNP Q82NB5
K	239	HIS	-	expression tag	UNP Q82NB5
K	240	HIS	-	expression tag	UNP Q82NB5
K	241	HIS	-	expression tag	UNP Q82NB5
K	242	HIS	-	expression tag	UNP Q82NB5
K	243	HIS	-	expression tag	UNP Q82NB5
L	18	MET	-	expression tag	UNP Q82NB5
L	19	SER	-	expression tag	UNP Q82NB5
L	20	LEU	-	expression tag	UNP Q82NB5
L	236	GLU	-	expression tag	UNP Q82NB5
L	237	GLY	-	expression tag	UNP Q82NB5
L	238	HIS	-	expression tag	UNP Q82NB5
L	239	HIS	-	expression tag	UNP Q82NB5
L	240	HIS	-	expression tag	UNP Q82NB5
L	241	HIS	-	expression tag	UNP Q82NB5
L	242	HIS	-	expression tag	UNP Q82NB5

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Chain	Residue	Modelled	Actual	Comment	Reference
L	243	HIS	-	expression tag	UNP Q82NB5

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	15	Total	O	0	0
			15	15		
3	B	8	Total	O	0	0
			8	8		
3	C	18	Total	O	0	0
			18	18		
3	D	19	Total	O	0	0
			19	19		
3	E	27	Total	O	0	0
			27	27		
3	F	9	Total	O	0	0
			9	9		
3	G	13	Total	O	0	0
			13	13		

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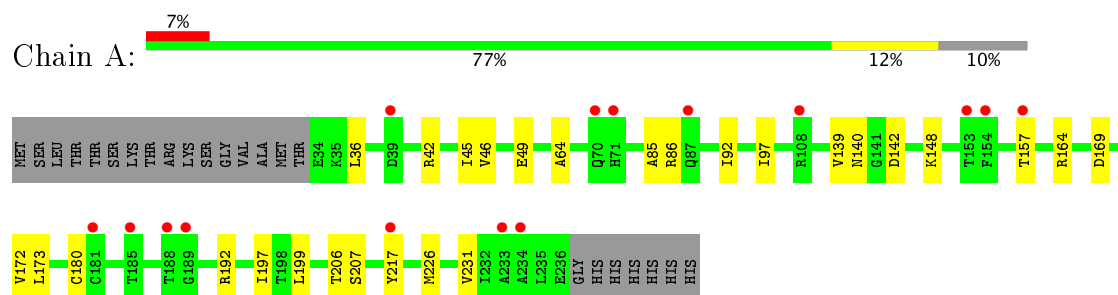
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	14	Total 14	O 14	0	0
3	I	12	Total 12	O 12	0	0
3	J	10	Total 10	O 10	0	0
3	K	22	Total 22	O 22	0	0
3	L	7	Total 7	O 7	0	0

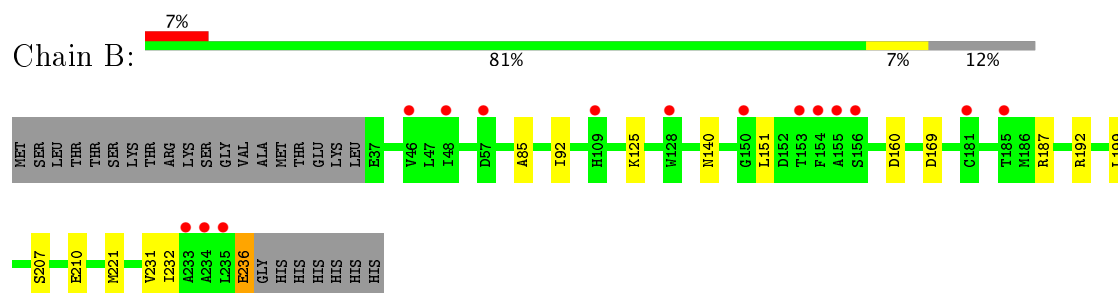
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

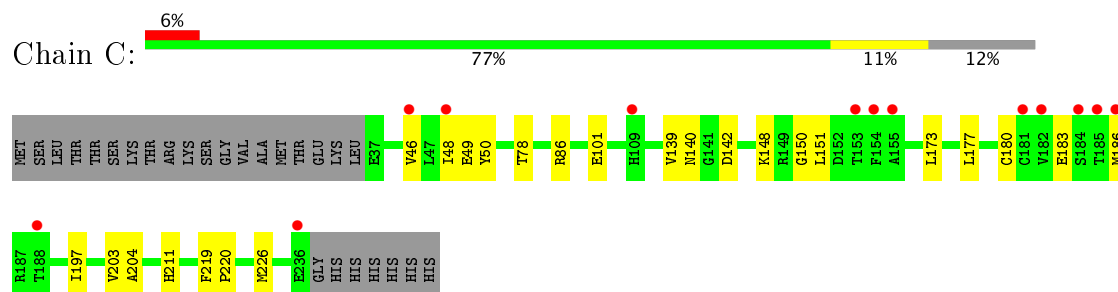
• Molecule 1: Putative isochorismatase



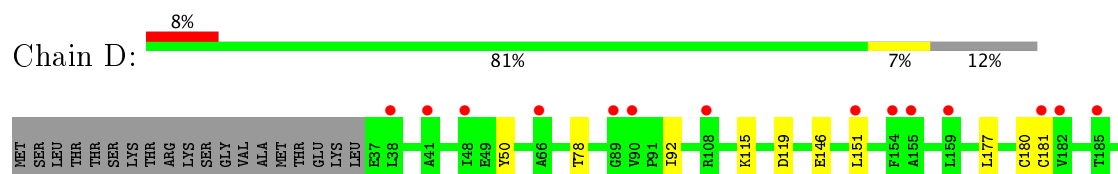
• Molecule 1: Putative isochorismatase

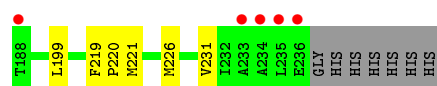


• Molecule 1: Putative isochorismatase

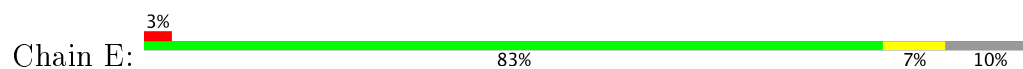


• Molecule 1: Putative isochorismatase

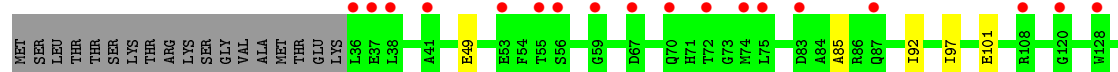
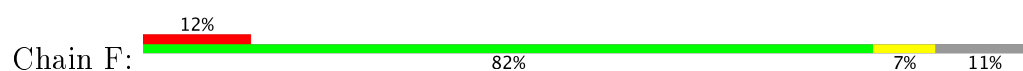




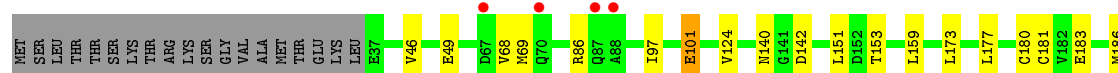
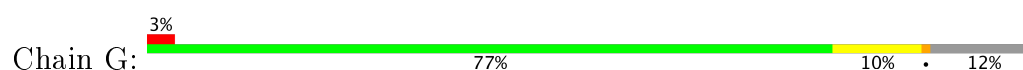
• Molecule 1: Putative isochorismatase



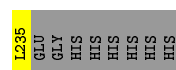
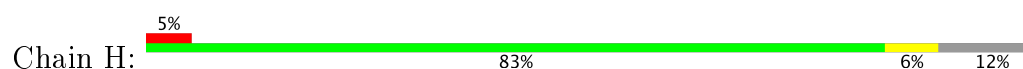
• Molecule 1: Putative isochorismatase



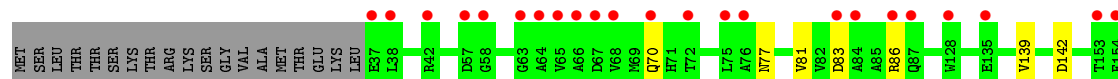
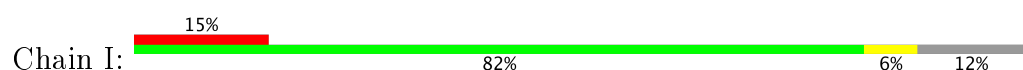
• Molecule 1: Putative isochorismatase

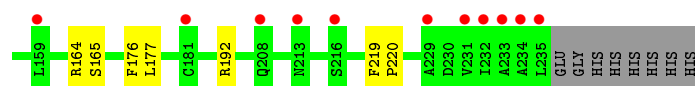


• Molecule 1: Putative isochorismatase

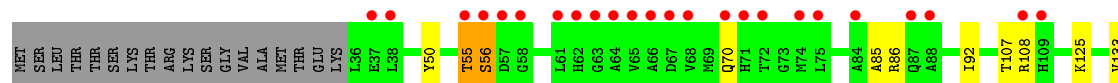
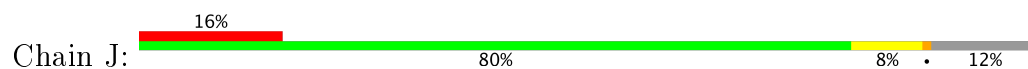


• Molecule 1: Putative isochorismatase

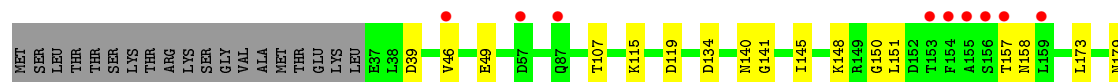
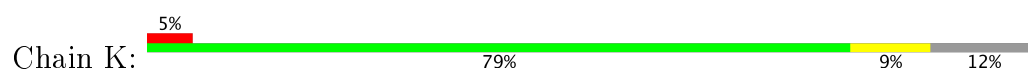




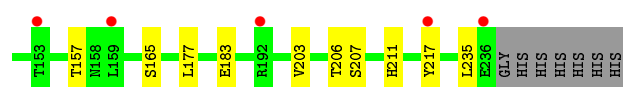
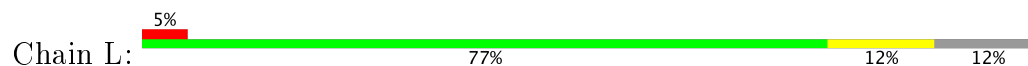
- Molecule 1: Putative isochorismatase



- Molecule 1: Putative isochorismatase



- Molecule 1: Putative isochorismatase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.74Å 112.10Å 178.90Å 90.00° 99.47° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 39.96 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.7 (20.00-2.30) 97.7 (39.96-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.79 (at 2.29Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.188 , 0.239 0.189 , 0.237	Depositor DCC
R_{free} test set	5723 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	31.8	Xtriage
Anisotropy	0.838	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 34.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.016 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17957	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.17% 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: 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	0.83	0/1518	0.79	0/2069
1	B	0.80	0/1487	0.76	1/2029 (0.0%)
1	C	0.86	0/1502	0.82	0/2046
1	D	0.82	0/1515	0.77	0/2062
1	E	0.86	0/1542	0.84	1/2099 (0.0%)
1	F	0.74	0/1520	0.73	0/2069
1	G	0.87	1/1515 (0.1%)	0.80	1/2062 (0.0%)
1	H	0.82	0/1512	0.79	0/2058
1	I	0.71	0/1481	0.69	0/2017
1	J	0.69	0/1495	0.73	0/2036
1	K	0.87	0/1502	0.79	0/2046
1	L	0.78	0/1507	0.78	1/2052 (0.0%)
All	All	0.80	1/18096 (0.0%)	0.77	4/24645 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	210	GLU	CG-CD	6.06	1.61	1.51

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	L	119	ASP	CB-CG-OD1	6.32	123.99	118.30
1	B	187	ARG	NE-CZ-NH1	5.84	123.22	120.30
1	E	164	ARG	NE-CZ-NH2	-5.43	117.58	120.30
1	G	69	MET	CG-SD-CE	-5.20	91.89	100.20

There are no chirality outliers.

There are no planarity outliers.

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	1491	0	1454	17	0
1	B	1460	0	1415	10	0
1	C	1475	0	1443	14	0
1	D	1488	0	1462	13	0
1	E	1513	0	1477	8	0
1	F	1493	0	1471	7	0
1	G	1488	0	1462	15	0
1	H	1485	0	1460	9	0
1	I	1456	0	1434	6	0
1	J	1469	0	1437	14	0
1	K	1475	0	1445	15	0
1	L	1480	0	1452	12	0
2	B	5	0	0	0	0
2	K	5	0	0	1	0
3	A	15	0	0	0	0
3	B	8	0	0	0	0
3	C	18	0	0	0	0
3	D	19	0	0	0	0
3	E	27	0	0	0	0
3	F	9	0	0	0	0
3	G	13	0	0	1	0
3	H	14	0	0	1	0
3	I	12	0	0	0	0
3	J	10	0	0	0	0
3	K	22	0	0	2	0
3	L	7	0	0	0	0
All	All	17957	0	17412	130	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 4.

All (130) 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:36:LEU:HD22	1:A:226:MET:SD	2.00	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:CYS:HB2	1:B:221:MET:HE1	1.61	0.83
1:C:197:ILE:HG12	1:C:226:MET:HE2	1.61	0.81
1:E:197:ILE:HG23	1:E:226:MET:HG3	1.62	0.81
1:J:55:THR:HG21	1:J:133:VAL:HB	1.61	0.81
1:G:183:GLU:HG2	1:G:186:MET:HE1	1.69	0.75
1:J:55:THR:HG21	1:J:133:VAL:CG2	2.17	0.74
1:D:226:MET:HE2	1:D:231:VAL:HG22	1.68	0.74
1:J:55:THR:HG21	1:J:133:VAL:CB	2.18	0.73
1:J:164:ARG:NH2	1:J:192:ARG:O	2.21	0.73
1:D:226:MET:CE	1:D:231:VAL:HG22	2.18	0.73
1:K:141:GLY:HA3	3:K:253:HOH:O	1.89	0.71
1:G:199:LEU:HD21	1:G:231:VAL:HG21	1.72	0.71
1:D:199:LEU:HD21	1:D:231:VAL:HG21	1.77	0.67
1:B:160:ASP:OD2	1:B:192:ARG:NH1	2.28	0.65
1:I:86:ARG:NH2	1:I:142:ASP:OD2	2.28	0.65
1:D:115:LYS:NZ	1:D:119:ASP:OD2	2.30	0.65
1:A:164:ARG:NH2	1:A:192:ARG:O	2.31	0.64
1:C:180:CYS:HB2	1:D:221:MET:CE	2.27	0.64
1:L:148:LYS:NZ	1:L:150:GLY:O	2.32	0.62
1:A:217:TYR:CD2	1:B:210:GLU:HG2	2.35	0.62
1:G:199:LEU:CD2	1:G:231:VAL:HG21	2.30	0.62
1:E:115:LYS:NZ	1:E:119:ASP:OD2	2.34	0.61
1:E:197:ILE:CG2	1:E:226:MET:HG3	2.30	0.60
1:C:46:VAL:HB	1:C:173:LEU:HD23	1.84	0.59
1:C:180:CYS:HB2	1:D:221:MET:HE1	1.84	0.59
1:K:179:ASN:HA	1:K:183:GLU:HG3	1.84	0.59
1:K:213:ASN:HD22	1:L:217:TYR:HE2	1.52	0.57
1:E:213:ASN:ND2	1:F:213:ASN:ND2	2.54	0.56
1:B:199:LEU:HD21	1:B:231:VAL:HG21	1.86	0.56
1:A:199:LEU:HD21	1:A:231:VAL:HG21	1.87	0.56
1:K:145:ILE:HG23	1:K:158:ASN:OD1	2.06	0.56
1:J:55:THR:HG21	1:J:133:VAL:HG21	1.88	0.56
1:C:151:LEU:HD22	1:C:151:LEU:N	2.22	0.55
1:D:92:ILE:HD13	1:D:92:ILE:N	2.22	0.54
1:G:101:GLU:HG2	1:G:124:VAL:CG1	2.37	0.54
1:J:85:ALA:HB3	1:J:92:ILE:HD11	1.89	0.54
1:G:101:GLU:HG2	1:G:124:VAL:HG11	1.90	0.53
1:I:164:ARG:NH2	1:I:192:ARG:O	2.41	0.53
1:A:85:ALA:HB3	1:A:92:ILE:HD11	1.91	0.53
1:K:145:ILE:HG22	1:K:157:THR:HB	1.91	0.53
1:G:49:GLU:HG2	1:G:97:ILE:HG22	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:177:LEU:HD22	1:J:205:ALA:HA	1.90	0.53
1:I:165:SER:O	1:K:107:THR:HB	2.09	0.53
1:A:45:ILE:HD13	1:A:172:VAL:HB	1.91	0.52
1:F:85:ALA:HB3	1:F:92:ILE:HD11	1.91	0.52
1:B:151:LEU:N	1:B:151:LEU:HD22	2.24	0.52
1:H:151:LEU:HD22	1:H:151:LEU:N	2.24	0.52
1:L:177:LEU:HD22	1:L:206:THR:HG23	1.92	0.52
1:L:203:VAL:HG23	1:L:211:HIS:HE1	1.74	0.51
1:A:86:ARG:NH2	1:A:142:ASP:OD2	2.43	0.51
1:B:199:LEU:CD2	1:B:231:VAL:HG21	2.41	0.51
1:H:50:TYR:CE1	1:H:78:THR:HG21	2.46	0.51
1:C:50:TYR:CZ	1:C:78:THR:HG21	2.47	0.50
1:G:151:LEU:HD12	1:G:181:CYS:SG	2.51	0.50
1:J:55:THR:HG22	1:J:56:SER:OG	2.11	0.50
1:D:151:LEU:HD12	1:D:181:CYS:SG	2.52	0.50
1:H:85:ALA:HA	1:H:235:LEU:HD13	1.93	0.50
1:F:151:LEU:HD22	1:F:151:LEU:N	2.28	0.49
1:A:199:LEU:CD2	1:A:231:VAL:HG21	2.43	0.49
1:L:90:VAL:HG21	1:L:235:LEU:CD2	2.43	0.49
1:A:85:ALA:CB	1:A:92:ILE:HD11	2.43	0.48
1:G:207:SER:HB3	1:G:210:GLU:OE1	2.13	0.48
1:I:77:ASN:O	1:I:81:VAL:HG23	2.12	0.48
1:A:197:ILE:HG23	1:A:226:MET:HG3	1.96	0.48
1:E:47:LEU:HD12	1:E:94:HIS:CE1	2.49	0.48
1:G:86:ARG:NH2	1:G:142:ASP:OD2	2.39	0.47
1:G:197:ILE:HG23	1:G:226:MET:HG3	1.96	0.47
1:J:107:THR:HB	1:L:165:SER:O	2.14	0.47
1:E:113:ILE:HD13	1:E:180:CYS:SG	2.55	0.47
1:G:153:THR:CG2	1:G:159:LEU:HD22	2.44	0.47
1:E:199:LEU:HD21	1:E:231:VAL:HG21	1.97	0.47
1:K:115:LYS:NZ	1:K:119:ASP:OD2	2.47	0.47
1:K:46:VAL:HB	1:K:173:LEU:HD23	1.97	0.46
1:A:49:GLU:HG2	1:A:97:ILE:HG22	1.97	0.46
1:G:177:LEU:HB2	1:G:180:CYS:HB3	1.98	0.46
1:D:50:TYR:CZ	1:D:78:THR:HG21	2.51	0.45
1:L:60:VAL:HG12	1:L:61:LEU:HG	1.98	0.45
1:D:199:LEU:CD2	1:D:231:VAL:HG21	2.43	0.45
1:G:68:VAL:HG23	3:G:4:HOH:O	2.15	0.45
1:K:205:ALA:HB3	1:K:211:HIS:HB2	1.98	0.45
1:C:86:ARG:NH2	1:C:142:ASP:OD2	2.48	0.45
1:K:148:LYS:NZ	1:K:150:GLY:O	2.44	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:148:LYS:NZ	1:C:150:GLY:O	2.41	0.45
1:C:183:GLU:HG3	1:C:186:MET:HE2	1.99	0.44
1:A:46:VAL:HB	1:A:173:LEU:HD23	1.99	0.44
1:B:85:ALA:HB3	1:B:92:ILE:HD11	1.98	0.44
1:G:46:VAL:HB	1:G:173:LEU:HD23	1.98	0.44
1:J:50:TYR:CD1	1:J:136:LEU:CD1	3.01	0.44
1:J:170:THR:HA	1:J:195:ARG:O	2.17	0.44
1:E:47:LEU:HB3	1:E:50:TYR:CD2	2.52	0.44
1:K:134:ASP:OD2	2:K:1:SO4:O2	2.35	0.44
1:L:43:THR:HG22	1:L:44:ALA:N	2.33	0.44
1:A:86:ARG:NH2	1:A:139:VAL:HG22	2.32	0.44
1:B:169:ASP:C	1:B:169:ASP:OD1	2.56	0.44
1:B:125:LYS:NZ	1:D:146:GLU:OE1	2.51	0.44
1:F:164:ARG:NH2	1:F:192:ARG:O	2.50	0.44
1:C:177:LEU:HA	1:C:204:ALA:O	2.19	0.43
1:L:133:VAL:HG22	1:L:135:GLU:OE1	2.19	0.43
1:K:39:ASP:C	1:K:39:ASP:OD2	2.56	0.43
1:C:203:VAL:HG23	1:C:211:HIS:HE1	1.84	0.43
1:C:219:PHE:N	1:C:220:PRO:CD	2.82	0.43
1:F:203:VAL:HG23	1:F:211:HIS:HE1	1.83	0.43
1:C:203:VAL:HG23	1:C:211:HIS:CE1	2.54	0.42
1:H:46:VAL:HB	1:H:173:LEU:HD23	2.01	0.42
1:H:155:ALA:HB3	3:H:2:HOH:O	2.20	0.42
1:A:169:ASP:C	1:A:169:ASP:OD1	2.58	0.42
1:H:151:LEU:HD22	1:H:151:LEU:H	1.84	0.42
1:J:176:PHE:HA	1:J:177:LEU:HA	1.80	0.42
1:D:177:LEU:HB2	1:D:180:CYS:HB3	2.01	0.42
1:H:148:LYS:NZ	1:H:150:GLY:O	2.47	0.42
1:K:140:ASN:O	3:K:253:HOH:O	2.22	0.42
1:F:160:ASP:O	1:F:161:PHE:C	2.59	0.42
1:G:217:TYR:CD2	1:H:210:GLU:HG2	2.54	0.42
1:K:49:GLU:OE1	1:K:148:LYS:NZ	2.49	0.42
1:B:232:ILE:O	1:B:236:GLU:HB2	2.19	0.41
1:F:49:GLU:HG2	1:F:97:ILE:HG22	2.03	0.41
1:K:151:LEU:HD22	1:K:151:LEU:N	2.35	0.41
1:C:48:ILE:HG22	1:C:49:GLU:HG3	2.02	0.41
1:H:85:ALA:HB1	1:H:90:VAL:HB	2.03	0.41
1:A:64:ALA:O	1:A:206:THR:HA	2.20	0.41
1:I:219:PHE:N	1:I:220:PRO:CD	2.83	0.41
1:D:219:PHE:N	1:D:220:PRO:CD	2.84	0.41
1:J:86:ARG:NH2	1:J:142:ASP:OD2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:145:ILE:HG22	1:L:157:THR:HB	2.03	0.41
1:I:176:PHE:HA	1:I:177:LEU:HA	1.89	0.41
1:L:47:LEU:HB3	1:L:50:TYR:CD2	2.56	0.41
1:J:86:ARG:HH21	1:J:142:ASP:CG	2.25	0.40
1:L:78:THR:O	1:L:82:VAL:HG23	2.21	0.40
1:A:148:LYS:N	1:A:157:THR:HG22	2.37	0.40

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	201/226 (89%)	197 (98%)	3 (2%)	1 (0%)	32	39
1	B	198/226 (88%)	195 (98%)	2 (1%)	1 (0%)	32	39
1	C	198/226 (88%)	193 (98%)	4 (2%)	1 (0%)	32	39
1	D	198/226 (88%)	194 (98%)	4 (2%)	0	100	100
1	E	202/226 (89%)	197 (98%)	4 (2%)	1 (0%)	32	39
1	F	199/226 (88%)	196 (98%)	2 (1%)	1 (0%)	32	39
1	G	198/226 (88%)	192 (97%)	5 (2%)	1 (0%)	32	39
1	H	198/226 (88%)	195 (98%)	3 (2%)	0	100	100
1	I	197/226 (87%)	191 (97%)	6 (3%)	0	100	100
1	J	198/226 (88%)	193 (98%)	5 (2%)	0	100	100
1	K	198/226 (88%)	197 (100%)	1 (0%)	0	100	100
1	L	198/226 (88%)	194 (98%)	4 (2%)	0	100	100
All	All	2383/2712 (88%)	2334 (98%)	43 (2%)	6 (0%)	44	55

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	140	ASN
1	G	140	ASN
1	A	140	ASN
1	B	140	ASN
1	E	140	ASN
1	F	140	ASN

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	153/181 (84%)	151 (99%)	2 (1%)	73	86
1	B	150/181 (83%)	148 (99%)	2 (1%)	73	86
1	C	153/181 (84%)	151 (99%)	2 (1%)	73	86
1	D	156/181 (86%)	156 (100%)	0	100	100
1	E	157/181 (87%)	155 (99%)	2 (1%)	73	86
1	F	156/181 (86%)	154 (99%)	2 (1%)	73	86
1	G	156/181 (86%)	154 (99%)	2 (1%)	73	86
1	H	155/181 (86%)	154 (99%)	1 (1%)	89	95
1	I	150/181 (83%)	147 (98%)	3 (2%)	60	77
1	J	151/181 (83%)	146 (97%)	5 (3%)	43	59
1	K	153/181 (84%)	153 (100%)	0	100	100
1	L	154/181 (85%)	149 (97%)	5 (3%)	44	60
All	All	1844/2172 (85%)	1818 (99%)	26 (1%)	71	85

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	ARG
1	A	207	SER
1	B	207	SER
1	B	236	GLU
1	C	101	GLU

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Mol	Chain	Res	Type
1	C	139	VAL
1	E	70	GLN
1	E	171	ILE
1	F	101	GLU
1	F	236	GLU
1	G	101	GLU
1	G	226	MET
1	H	101	GLU
1	I	70	GLN
1	I	83	ASP
1	I	139	VAL
1	J	55	THR
1	J	56	SER
1	J	70	GLN
1	J	108	ARG
1	J	125	LYS
1	L	42	ARG
1	L	60	VAL
1	L	139	VAL
1	L	183	GLU
1	L	207	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	208	GLN
1	C	87	GLN
1	D	87	GLN
1	D	94	HIS
1	E	208	GLN
1	E	213	ASN
1	F	109	HIS
1	F	213	ASN
1	G	208	GLN
1	H	208	GLN
1	I	70	GLN
1	K	208	GLN
1	K	213	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

2 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	2	-	4,4,4	0.29	0	6,6,6	0.45	0
2	SO4	K	1	-	4,4,4	0.25	0	6,6,6	0.47	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	SO4	B	2	-	-	0/0/0/0	0/0/0/0
2	SO4	K	1	-	-	0/0/0/0	0/0/0/0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	K	1	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 ⓘ

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	203/226 (89%)	0.59	15 (7%)	15 21	11, 22, 34, 50	0
1	B	200/226 (88%)	0.48	15 (7%)	15 20	12, 21, 31, 41	0
1	C	200/226 (88%)	0.41	13 (6%)	20 26	13, 20, 30, 38	0
1	D	200/226 (88%)	0.63	19 (9%)	9 12	12, 19, 30, 35	0
1	E	204/226 (90%)	0.26	6 (2%)	52 59	12, 20, 28, 32	0
1	F	201/226 (88%)	0.80	28 (13%)	3 5	12, 21, 30, 40	0
1	G	200/226 (88%)	0.33	6 (3%)	51 58	10, 21, 31, 39	0
1	H	200/226 (88%)	0.25	11 (5%)	26 32	11, 20, 30, 41	0
1	I	199/226 (88%)	0.83	34 (17%)	2 2	13, 22, 29, 34	0
1	J	200/226 (88%)	0.89	36 (18%)	2 2	12, 21, 31, 43	0
1	K	200/226 (88%)	0.37	11 (5%)	26 32	11, 19, 27, 33	0
1	L	200/226 (88%)	0.43	11 (5%)	26 32	14, 21, 29, 37	0
All	All	2407/2712 (88%)	0.52	205 (8%)	11 16	10, 20, 30, 50	0

All (205) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	66	ALA	5.8
1	J	217	TYR	5.0
1	J	70	GLN	4.8
1	F	83	ASP	4.5
1	I	75	LEU	4.5
1	C	181	CYS	4.3
1	H	36	LEU	4.2
1	I	84	ALA	4.2
1	J	75	LEU	4.2
1	D	235	LEU	4.1
1	J	58	GLY	4.1

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Mol	Chain	Res	Type	RSRZ
1	E	239	HIS	4.1
1	L	58	GLY	4.0
1	F	235	LEU	4.0
1	J	209	GLU	4.0
1	I	76	ALA	4.0
1	E	238	HIS	3.8
1	D	38	LEU	3.7
1	I	235	LEU	3.7
1	D	236	GLU	3.7
1	D	181	CYS	3.6
1	D	89	GLY	3.6
1	H	217	TYR	3.6
1	J	84	ALA	3.5
1	B	181	CYS	3.5
1	F	72	THR	3.5
1	J	109	HIS	3.5
1	I	87	GLN	3.5
1	J	56	SER	3.5
1	I	83	ASP	3.5
1	J	64	ALA	3.4
1	F	87	GLN	3.4
1	H	58	GLY	3.4
1	L	121	LYS	3.3
1	C	236	GLU	3.3
1	J	88	ALA	3.3
1	F	36	LEU	3.3
1	K	87	GLN	3.2
1	F	236	GLU	3.2
1	J	57	ASP	3.2
1	B	185	THR	3.2
1	J	61	LEU	3.2
1	G	87	GLN	3.2
1	H	157	THR	3.1
1	A	71	HIS	3.1
1	J	68	VAL	3.1
1	C	153	THR	3.1
1	A	70	GLN	3.1
1	C	185	THR	3.0
1	G	236	GLU	3.0
1	I	213	ASN	3.0
1	J	87	GLN	3.0
1	I	57	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
1	F	108	ARG	3.0
1	B	155	ALA	2.9
1	I	135	GLU	2.9
1	A	185	THR	2.9
1	L	217	TYR	2.9
1	H	181	CYS	2.9
1	F	128	TRP	2.9
1	D	155	ALA	2.8
1	K	159	LEU	2.8
1	I	38	LEU	2.8
1	I	86	ARG	2.8
1	H	233	ALA	2.8
1	B	153	THR	2.8
1	J	65	VAL	2.8
1	E	217	TYR	2.8
1	J	216	SER	2.8
1	D	154	PHE	2.7
1	C	184	SER	2.7
1	F	59	GLY	2.7
1	F	75	LEU	2.7
1	F	217	TYR	2.7
1	J	230	ASP	2.7
1	J	37	GLU	2.7
1	H	159	LEU	2.7
1	E	237	GLY	2.7
1	I	208	GLN	2.7
1	L	70	GLN	2.7
1	I	154	PHE	2.7
1	D	188	THR	2.7
1	I	232	ILE	2.7
1	I	66	ALA	2.7
1	G	217	TYR	2.7
1	K	57	ASP	2.7
1	K	46	VAL	2.7
1	J	232	ILE	2.7
1	B	154	PHE	2.7
1	I	65	VAL	2.7
1	B	234	ALA	2.6
1	A	153	THR	2.6
1	I	58	GLY	2.6
1	K	181	CYS	2.6
1	A	217	TYR	2.6

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Mol	Chain	Res	Type	RSRZ
1	J	38	LEU	2.6
1	I	70	GLN	2.6
1	I	67	ASP	2.6
1	I	229	ALA	2.6
1	G	88	ALA	2.6
1	I	216	SER	2.5
1	F	74	MET	2.5
1	L	236	GLU	2.5
1	J	233	ALA	2.5
1	B	235	LEU	2.5
1	J	62	HIS	2.5
1	A	154	PHE	2.5
1	D	41	ALA	2.5
1	D	159	LEU	2.5
1	I	159	LEU	2.5
1	K	154	PHE	2.5
1	D	233	ALA	2.5
1	F	41	ALA	2.5
1	C	48	ILE	2.5
1	L	38	LEU	2.5
1	B	46	VAL	2.5
1	K	156	SER	2.5
1	J	153	THR	2.5
1	L	40	PRO	2.5
1	H	234	ALA	2.4
1	B	48	ILE	2.4
1	J	108	ARG	2.4
1	F	37	GLU	2.4
1	C	154	PHE	2.4
1	I	153	THR	2.4
1	F	67	ASP	2.4
1	E	181	CYS	2.4
1	I	72	THR	2.4
1	J	55	THR	2.4
1	J	72	THR	2.4
1	B	128	TRP	2.4
1	F	159	LEU	2.4
1	J	63	GLY	2.4
1	K	155	ALA	2.4
1	B	156	SER	2.4
1	F	53	GLU	2.4
1	K	157	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	90	VAL	2.4
1	E	154	PHE	2.4
1	I	181	CYS	2.4
1	B	57	ASP	2.3
1	G	67	ASP	2.3
1	J	67	ASP	2.3
1	D	48	ILE	2.3
1	I	37	GLU	2.3
1	D	185	THR	2.3
1	I	42	ARG	2.3
1	J	74	MET	2.3
1	F	231	VAL	2.3
1	B	109	HIS	2.3
1	D	66	ALA	2.3
1	F	38	LEU	2.3
1	L	159	LEU	2.3
1	K	153	THR	2.3
1	F	154	PHE	2.3
1	I	63	GLY	2.3
1	A	181	CYS	2.3
1	B	233	ALA	2.3
1	H	41	ALA	2.3
1	J	159	LEU	2.3
1	A	189	GLY	2.3
1	C	46	VAL	2.3
1	I	68	VAL	2.3
1	A	233	ALA	2.3
1	A	108	ARG	2.2
1	A	188	THR	2.2
1	L	192	ARG	2.2
1	J	212	ASN	2.2
1	D	108	ARG	2.2
1	A	157	THR	2.2
1	C	188	THR	2.2
1	F	55	THR	2.2
1	I	234	ALA	2.2
1	D	182	VAL	2.2
1	A	39	ASP	2.2
1	F	56	SER	2.2
1	F	70	GLN	2.2
1	J	71	HIS	2.2
1	I	128	TRP	2.2

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Mol	Chain	Res	Type	RSRZ
1	F	201	ASP	2.2
1	A	87	GLN	2.2
1	C	182	VAL	2.2
1	F	213	ASN	2.2
1	H	154	PHE	2.2
1	D	234	ALA	2.2
1	C	186	MET	2.2
1	C	109	HIS	2.1
1	D	151	LEU	2.1
1	H	141	GLY	2.1
1	K	185	THR	2.1
1	J	231	VAL	2.1
1	L	153	THR	2.1
1	L	42	ARG	2.1
1	F	120	GLY	2.1
1	I	233	ALA	2.1
1	G	70	GLN	2.1
1	A	234	ALA	2.1
1	I	64	ALA	2.1
1	F	133	VAL	2.1
1	F	233	ALA	2.1
1	J	155	ALA	2.1
1	I	231	VAL	2.0
1	C	155	ALA	2.0
1	J	211	HIS	2.0
1	B	150	GLY	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	SO4	K	1	5/5	0.98	0.14	-0.08	31,33,34,36	0
2	SO4	B	2	5/5	0.94	0.19	-0.51	63,65,66,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.