



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:41 PM GMT

PDB ID : 3ROJ
Title : D-fructose 1,6-bisphosphatase class 2/sedoheptulose 1,7-bisphosphatase of Synechocystis sp. PCC 6803
Authors : Hu, X.; Hui, D.; Lingling, F.; Jian, W.
Deposited on : 2011-04-26
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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

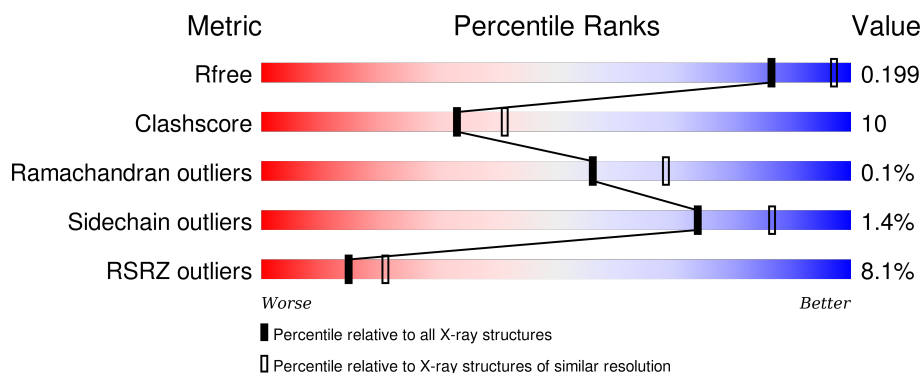
1 Overall quality at a glance

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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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	379	 7% 75% 16% 8%
1	B	379	 8% 72% 18% 9%
1	C	379	 7% 75% 16% 8%
1	D	379	 7% 73% 18% 9%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MG	A	347	-	-	-	X
2	MG	B	347	-	-	-	X
2	MG	C	347	-	-	-	X
4	SO4	A	350	-	-	-	X
4	SO4	C	352	-	-	-	X
4	SO4	D	350	-	-	-	X
5	CL	A	355	-	-	X	X
5	CL	A	358	-	-	-	X
5	CL	A	361	-	-	X	-
5	CL	B	351	-	-	X	-
6	GOL	A	365	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 11412 atoms, of which 8 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 D-fructose 1,6-bisphosphatase class 2/sedoheptulose 1,7-bisphosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	347	Total	C	N	O	S	0	4	0
			2591	1607	455	507	22			
1	B	345	Total	C	N	O	S	0	4	0
			2576	1599	454	501	22			
1	C	347	Total	C	N	O	S	0	2	0
			2563	1590	452	499	22			
1	D	346	Total	C	N	O	S	0	2	0
			2579	1602	453	502	22			

There are 140 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-33	MET	-	EXPRESSION TAG	UNP P73922
A	-32	GLY	-	EXPRESSION TAG	UNP P73922
A	-31	SER	-	EXPRESSION TAG	UNP P73922
A	-30	SER	-	EXPRESSION TAG	UNP P73922
A	-29	HIS	-	EXPRESSION TAG	UNP P73922
A	-28	HIS	-	EXPRESSION TAG	UNP P73922
A	-27	HIS	-	EXPRESSION TAG	UNP P73922
A	-26	HIS	-	EXPRESSION TAG	UNP P73922
A	-25	HIS	-	EXPRESSION TAG	UNP P73922
A	-24	HIS	-	EXPRESSION TAG	UNP P73922
A	-23	SER	-	EXPRESSION TAG	UNP P73922
A	-22	SER	-	EXPRESSION TAG	UNP P73922
A	-21	GLY	-	EXPRESSION TAG	UNP P73922
A	-20	LEU	-	EXPRESSION TAG	UNP P73922
A	-19	VAL	-	EXPRESSION TAG	UNP P73922
A	-18	PRO	-	EXPRESSION TAG	UNP P73922
A	-17	ARG	-	EXPRESSION TAG	UNP P73922
A	-16	GLY	-	EXPRESSION TAG	UNP P73922
A	-15	SER	-	EXPRESSION TAG	UNP P73922
A	-14	HIS	-	EXPRESSION TAG	UNP P73922
A	-13	MET	-	EXPRESSION TAG	UNP P73922

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-12	ALA	-	EXPRESSION TAG	UNP P73922
A	-11	SER	-	EXPRESSION TAG	UNP P73922
A	-10	MET	-	EXPRESSION TAG	UNP P73922
A	-9	THR	-	EXPRESSION TAG	UNP P73922
A	-8	GLY	-	EXPRESSION TAG	UNP P73922
A	-7	GLY	-	EXPRESSION TAG	UNP P73922
A	-6	GLN	-	EXPRESSION TAG	UNP P73922
A	-5	GLN	-	EXPRESSION TAG	UNP P73922
A	-4	MET	-	EXPRESSION TAG	UNP P73922
A	-3	GLY	-	EXPRESSION TAG	UNP P73922
A	-2	ARG	-	EXPRESSION TAG	UNP P73922
A	-1	GLY	-	EXPRESSION TAG	UNP P73922
A	0	SER	-	EXPRESSION TAG	UNP P73922
A	1	VAL	-	EXPRESSION TAG	UNP P73922
B	-33	MET	-	EXPRESSION TAG	UNP P73922
B	-32	GLY	-	EXPRESSION TAG	UNP P73922
B	-31	SER	-	EXPRESSION TAG	UNP P73922
B	-30	SER	-	EXPRESSION TAG	UNP P73922
B	-29	HIS	-	EXPRESSION TAG	UNP P73922
B	-28	HIS	-	EXPRESSION TAG	UNP P73922
B	-27	HIS	-	EXPRESSION TAG	UNP P73922
B	-26	HIS	-	EXPRESSION TAG	UNP P73922
B	-25	HIS	-	EXPRESSION TAG	UNP P73922
B	-24	HIS	-	EXPRESSION TAG	UNP P73922
B	-23	SER	-	EXPRESSION TAG	UNP P73922
B	-22	SER	-	EXPRESSION TAG	UNP P73922
B	-21	GLY	-	EXPRESSION TAG	UNP P73922
B	-20	LEU	-	EXPRESSION TAG	UNP P73922
B	-19	VAL	-	EXPRESSION TAG	UNP P73922
B	-18	PRO	-	EXPRESSION TAG	UNP P73922
B	-17	ARG	-	EXPRESSION TAG	UNP P73922
B	-16	GLY	-	EXPRESSION TAG	UNP P73922
B	-15	SER	-	EXPRESSION TAG	UNP P73922
B	-14	HIS	-	EXPRESSION TAG	UNP P73922
B	-13	MET	-	EXPRESSION TAG	UNP P73922
B	-12	ALA	-	EXPRESSION TAG	UNP P73922
B	-11	SER	-	EXPRESSION TAG	UNP P73922
B	-10	MET	-	EXPRESSION TAG	UNP P73922
B	-9	THR	-	EXPRESSION TAG	UNP P73922
B	-8	GLY	-	EXPRESSION TAG	UNP P73922
B	-7	GLY	-	EXPRESSION TAG	UNP P73922
B	-6	GLN	-	EXPRESSION TAG	UNP P73922

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-5	GLN	-	EXPRESSION TAG	UNP P73922
B	-4	MET	-	EXPRESSION TAG	UNP P73922
B	-3	GLY	-	EXPRESSION TAG	UNP P73922
B	-2	ARG	-	EXPRESSION TAG	UNP P73922
B	-1	GLY	-	EXPRESSION TAG	UNP P73922
B	0	SER	-	EXPRESSION TAG	UNP P73922
B	1	VAL	-	EXPRESSION TAG	UNP P73922
C	-33	MET	-	EXPRESSION TAG	UNP P73922
C	-32	GLY	-	EXPRESSION TAG	UNP P73922
C	-31	SER	-	EXPRESSION TAG	UNP P73922
C	-30	SER	-	EXPRESSION TAG	UNP P73922
C	-29	HIS	-	EXPRESSION TAG	UNP P73922
C	-28	HIS	-	EXPRESSION TAG	UNP P73922
C	-27	HIS	-	EXPRESSION TAG	UNP P73922
C	-26	HIS	-	EXPRESSION TAG	UNP P73922
C	-25	HIS	-	EXPRESSION TAG	UNP P73922
C	-24	HIS	-	EXPRESSION TAG	UNP P73922
C	-23	SER	-	EXPRESSION TAG	UNP P73922
C	-22	SER	-	EXPRESSION TAG	UNP P73922
C	-21	GLY	-	EXPRESSION TAG	UNP P73922
C	-20	LEU	-	EXPRESSION TAG	UNP P73922
C	-19	VAL	-	EXPRESSION TAG	UNP P73922
C	-18	PRO	-	EXPRESSION TAG	UNP P73922
C	-17	ARG	-	EXPRESSION TAG	UNP P73922
C	-16	GLY	-	EXPRESSION TAG	UNP P73922
C	-15	SER	-	EXPRESSION TAG	UNP P73922
C	-14	HIS	-	EXPRESSION TAG	UNP P73922
C	-13	MET	-	EXPRESSION TAG	UNP P73922
C	-12	ALA	-	EXPRESSION TAG	UNP P73922
C	-11	SER	-	EXPRESSION TAG	UNP P73922
C	-10	MET	-	EXPRESSION TAG	UNP P73922
C	-9	THR	-	EXPRESSION TAG	UNP P73922
C	-8	GLY	-	EXPRESSION TAG	UNP P73922
C	-7	GLY	-	EXPRESSION TAG	UNP P73922
C	-6	GLN	-	EXPRESSION TAG	UNP P73922
C	-5	GLN	-	EXPRESSION TAG	UNP P73922
C	-4	MET	-	EXPRESSION TAG	UNP P73922
C	-3	GLY	-	EXPRESSION TAG	UNP P73922
C	-2	ARG	-	EXPRESSION TAG	UNP P73922
C	-1	GLY	-	EXPRESSION TAG	UNP P73922
C	0	SER	-	EXPRESSION TAG	UNP P73922
C	1	VAL	-	EXPRESSION TAG	UNP P73922

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-33	MET	-	EXPRESSION TAG	UNP P73922
D	-32	GLY	-	EXPRESSION TAG	UNP P73922
D	-31	SER	-	EXPRESSION TAG	UNP P73922
D	-30	SER	-	EXPRESSION TAG	UNP P73922
D	-29	HIS	-	EXPRESSION TAG	UNP P73922
D	-28	HIS	-	EXPRESSION TAG	UNP P73922
D	-27	HIS	-	EXPRESSION TAG	UNP P73922
D	-26	HIS	-	EXPRESSION TAG	UNP P73922
D	-25	HIS	-	EXPRESSION TAG	UNP P73922
D	-24	HIS	-	EXPRESSION TAG	UNP P73922
D	-23	SER	-	EXPRESSION TAG	UNP P73922
D	-22	SER	-	EXPRESSION TAG	UNP P73922
D	-21	GLY	-	EXPRESSION TAG	UNP P73922
D	-20	LEU	-	EXPRESSION TAG	UNP P73922
D	-19	VAL	-	EXPRESSION TAG	UNP P73922
D	-18	PRO	-	EXPRESSION TAG	UNP P73922
D	-17	ARG	-	EXPRESSION TAG	UNP P73922
D	-16	GLY	-	EXPRESSION TAG	UNP P73922
D	-15	SER	-	EXPRESSION TAG	UNP P73922
D	-14	HIS	-	EXPRESSION TAG	UNP P73922
D	-13	MET	-	EXPRESSION TAG	UNP P73922
D	-12	ALA	-	EXPRESSION TAG	UNP P73922
D	-11	SER	-	EXPRESSION TAG	UNP P73922
D	-10	MET	-	EXPRESSION TAG	UNP P73922
D	-9	THR	-	EXPRESSION TAG	UNP P73922
D	-8	GLY	-	EXPRESSION TAG	UNP P73922
D	-7	GLY	-	EXPRESSION TAG	UNP P73922
D	-6	GLN	-	EXPRESSION TAG	UNP P73922
D	-5	GLN	-	EXPRESSION TAG	UNP P73922
D	-4	MET	-	EXPRESSION TAG	UNP P73922
D	-3	GLY	-	EXPRESSION TAG	UNP P73922
D	-2	ARG	-	EXPRESSION TAG	UNP P73922
D	-1	GLY	-	EXPRESSION TAG	UNP P73922
D	0	SER	-	EXPRESSION TAG	UNP P73922
D	1	VAL	-	EXPRESSION TAG	UNP P73922

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

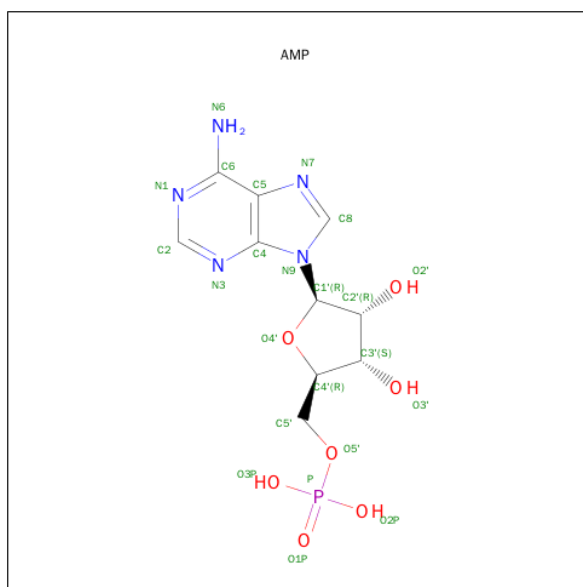
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	2	Total Mg 2 2	0	0
2	A	2	Total Mg 2 2	0	0

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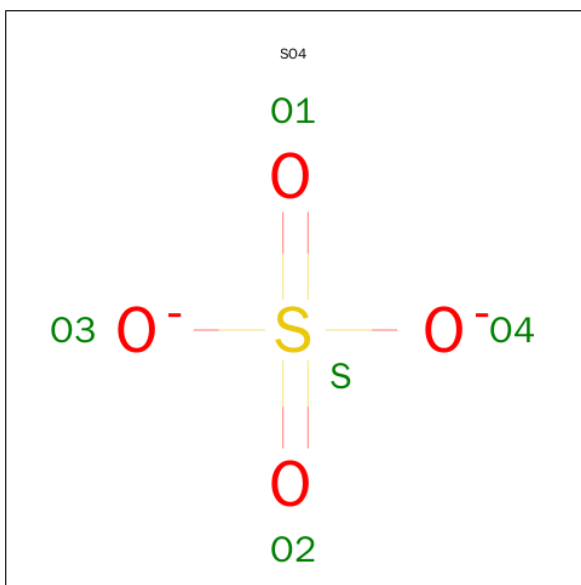
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	2	Total	Mg	0	0
			2	2		
2	C	2	Total	Mg	0	0
			2	2		

- Molecule 3 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula: $C_{10}H_{14}N_5O_7P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
3	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
3	C	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
3	C	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

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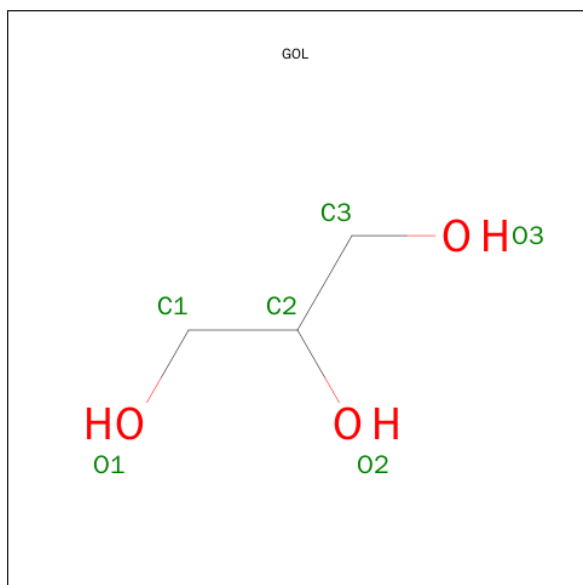
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	5	Total	Cl	0	0
			5	5		
5	A	10	Total	Cl	0	0
			10	10		
5	D	2	Total	Cl	0	0
			2	2		
5	C	6	Total	Cl	0	0
			6	6		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



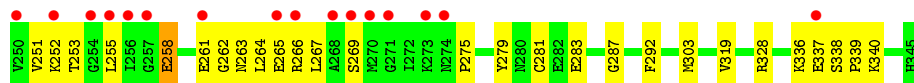
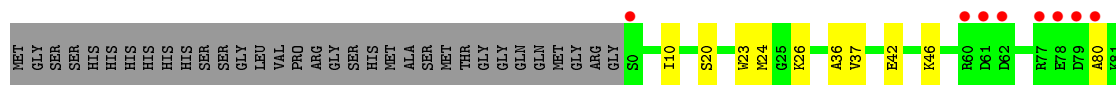
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	239	Total 239	O 239	0	0
7	B	209	Total 209	O 209	0	0
7	C	237	Total 237	O 237	0	0
7	D	196	Total 196	O 196	0	0



- Molecule 1: D-fructose 1,6-bisphosphatase class 2/sedoheptulose 1,7-bisphosphatase



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	144.28Å 144.28Å 168.79Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.54 – 2.30 25.54 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.3 (25.54-2.30) 99.3 (25.54-2.30)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 2.31Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.183 , 0.204 0.173 , 0.199	Depositor DCC
R_{free} test set	4393 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.392	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 55.1	EDS
Estimated twinning fraction	0.024 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 87538 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11412	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, MG, CL, CSO, SO4, AMP

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	0.25	0/2630	0.45	0/3550
1	B	0.25	0/2615	0.43	0/3532
1	C	0.26	0/2596	0.44	0/3507
1	D	0.24	0/2612	0.42	0/3525
All	All	0.25	0/10453	0.43	0/14114

There are no bond length outliers.

There are no bond angle outliers.

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	2591	0	2564	54	0
1	B	2576	0	2552	64	0
1	C	2563	0	2525	46	0
1	D	2579	0	2563	48	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	23	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	23	0	12	0	0
3	C	46	0	24	1	0
4	A	30	0	0	0	0
4	B	10	0	0	0	0
4	C	20	0	0	2	0
4	D	25	0	0	1	0
5	A	10	0	0	7	0
5	B	5	0	0	3	0
5	C	6	0	0	0	0
5	D	2	0	0	1	0
6	A	6	8	8	2	0
7	A	239	0	0	4	0
7	B	209	0	0	5	0
7	C	237	0	0	5	0
7	D	196	0	0	0	0
All	All	11404	8	10260	213	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 10.

All (213) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:263:ASN:HA	1:D:266:ARG:HD2	1.37	1.07
1:B:301:THR:O	7:B:707:HOH:O	1.92	0.85
1:A:150:SER:OG	1:A:153[A]:GLU:HG3	1.78	0.84
1:D:253:THR:HG22	1:D:258:GLU:CD	2.00	0.82
1:B:277[B]:GLN:HE21	1:B:277[B]:GLN:HA	1.49	0.78
1:B:113:MET:HE2	1:B:303:MET:CE	2.13	0.77
1:A:113:MET:HE2	1:A:303:MET:CE	2.15	0.76
1:A:280:ASN:HB2	6:A:365:GOL:H2	1.68	0.76
1:A:273:LYS:HG2	1:A:274:ASN:H	1.50	0.76
1:D:150:SER:OG	1:D:153[B]:GLU:HG2	1.84	0.75
1:A:273:LYS:CD	1:A:273:LYS:H	1.99	0.75
1:C:266:ARG:O	1:C:270:MET:HG3	1.87	0.74
1:B:81:LYS:HA	1:B:89:LEU:HD11	1.70	0.74
1:A:234:ARG:NE	5:A:355:CL:CL	2.56	0.73
1:A:273:LYS:HD3	1:A:273:LYS:H	1.53	0.73
1:A:273:LYS:HD3	1:A:273:LYS:N	2.03	0.73
1:D:262:GLY:O	1:D:266:ARG:HG3	1.90	0.72
1:D:85:ASN:HB2	1:D:86:PRO:CD	2.20	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:MET:CE	1:B:303:MET:CE	2.68	0.71
1:C:113:MET:HE2	1:C:303:MET:CE	2.20	0.71
1:B:113:MET:CE	1:B:303:MET:HE1	2.22	0.70
1:B:277[B]:GLN:NE2	1:B:277[B]:GLN:HA	2.05	0.70
5:B:351:CL:CL	7:B:600:HOH:O	2.47	0.70
1:D:253:THR:HG22	1:D:258:GLU:OE2	1.90	0.69
1:D:85:ASN:HB2	1:D:86:PRO:HD2	1.75	0.69
1:B:90:VAL:O	7:B:683:HOH:O	2.09	0.69
1:A:113:MET:HE2	1:A:303:MET:HE3	1.75	0.69
1:B:85:ASN:HB2	1:B:86:PRO:HD2	1.75	0.68
1:B:277[B]:GLN:NE2	1:B:278:VAL:H	1.91	0.68
1:C:194:ARG:NE	7:C:732:HOH:O	2.28	0.66
1:B:262:GLY:O	1:B:266:ARG:HG3	1.95	0.66
1:D:138:PRO:HD3	1:D:215:HIS:O	1.96	0.66
1:A:274:ASN:ND2	1:A:277[B]:GLN:HG2	2.10	0.65
1:A:113:MET:CE	1:A:303:MET:CE	2.74	0.65
1:A:136:ALA:HB3	1:A:217:LEU:HB3	1.79	0.64
1:C:136:ALA:HB3	1:C:217:LEU:HB3	1.79	0.64
5:B:351:CL:CL	7:B:538:HOH:O	2.52	0.64
1:B:80:ALA:CB	1:B:89:LEU:HD21	2.28	0.64
1:C:113:MET:CE	1:C:303:MET:CE	2.77	0.63
1:B:113:MET:HE2	1:B:303:MET:HE1	1.78	0.63
1:C:336:LYS:HA	7:C:781:HOH:O	1.99	0.63
1:B:23:TRP:CE3	1:B:26:LYS:HD3	2.33	0.62
1:B:128:PRO:HB3	1:B:270:MET:HE3	1.81	0.62
1:D:97:ASP:OD1	1:D:100:GLU:HG2	2.00	0.62
1:A:273:LYS:O	1:A:275:PRO:HD3	2.00	0.62
1:D:136:ALA:HB3	1:D:217:LEU:HB3	1.83	0.60
1:A:311:GLY:O	1:A:339:PRO:HG3	2.01	0.60
1:B:97:ASP:HB3	1:B:115:VAL:HG13	1.84	0.60
1:D:253:THR:HG23	1:D:255:LEU:H	1.67	0.60
1:B:263:ASN:HA	1:B:266:ARG:HD2	1.84	0.60
1:B:164:ARG:HD2	1:B:169:LEU:HD23	1.85	0.59
1:B:338:SER:N	1:B:339:PRO:CD	2.65	0.59
5:A:358:CL:CL	7:A:577:HOH:O	2.54	0.59
1:C:97:ASP:OD1	1:C:100:GLU:HG2	2.02	0.59
1:B:248:PRO:HA	1:B:251:VAL:O	2.03	0.59
1:A:280:ASN:CB	6:A:365:GOL:H2	2.32	0.58
1:D:336:LYS:HA	1:D:337:GLU:HB2	1.86	0.58
1:A:101:GLY:N	5:A:361:CL:CL	2.72	0.58
1:C:38:GLU:HG2	1:C:65:MET:CE	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:273:LYS:N	1:A:273:LYS:CD	2.62	0.57
1:D:338:SER:N	1:D:339:PRO:CD	2.68	0.57
1:B:23:TRP:CZ3	1:B:26:LYS:HD3	2.40	0.57
1:D:261:GLU:O	1:D:265:GLU:HG3	2.05	0.56
1:B:142:LYS:HA	1:B:239:HIS:HB2	1.88	0.56
1:A:273:LYS:HG2	1:A:274:ASN:N	2.20	0.56
1:B:97:ASP:HB3	1:B:115:VAL:CG1	2.35	0.56
1:C:10:ILE:HG12	1:C:319:VAL:HG21	1.87	0.56
1:C:205:ILE:HG22	1:C:302:LEU:HD11	1.86	0.56
1:D:80:ALA:HA	1:D:83:PHE:CE1	2.41	0.56
1:A:20:SER:HB2	1:A:36:ALA:HB2	1.88	0.55
1:C:336:LYS:CA	7:C:781:HOH:O	2.54	0.55
1:B:184:GLN:NE2	7:B:694:HOH:O	2.41	0.54
1:C:142:LYS:HA	1:C:239:HIS:HB2	1.89	0.54
1:D:130:PHE:CE2	1:D:267:LEU:HD21	2.43	0.53
1:C:38:GLU:HG2	1:C:65:MET:SD	2.48	0.53
1:C:20:SER:HB2	1:C:36:ALA:HB2	1.89	0.53
1:B:9:ILE:HD12	1:B:293:ALA:HB2	1.91	0.53
1:C:85:ASN:HB2	1:C:86:PRO:HD2	1.89	0.53
1:B:80:ALA:HB1	1:B:84:CYS:SG	2.49	0.52
1:B:80:ALA:HB3	1:B:89:LEU:HD21	1.90	0.52
1:B:37:VAL:HG22	1:B:98:PRO:HG2	1.91	0.52
1:B:20:SER:HB2	1:B:36:ALA:HB2	1.92	0.52
1:D:240:PHE:HB3	1:D:281:CYS:HB3	1.92	0.52
1:C:85:ASN:HB2	1:C:86:PRO:CD	2.41	0.51
1:B:85:ASN:HB2	1:B:86:PRO:CD	2.39	0.51
1:C:37:VAL:HG22	1:C:98:PRO:HG2	1.93	0.51
1:B:113:MET:HE1	1:B:303:MET:HE1	1.93	0.51
1:B:116:LEU:HD12	1:B:116:LEU:C	2.31	0.50
1:D:37:VAL:HG22	1:D:98:PRO:HG2	1.93	0.50
1:D:42:GLU:OE2	1:D:46:LYS:HE3	2.11	0.50
1:A:277[B]:GLN:NE2	7:A:591:HOH:O	2.43	0.50
1:C:24:MET:HG3	1:C:108:GLY:HA2	1.92	0.50
1:D:119:SER:OG	1:D:287:GLY:HA3	2.12	0.50
1:D:113:MET:SD	1:D:303:MET:CE	3.00	0.50
1:C:253:THR:HB	1:C:258:GLU:OE2	2.12	0.50
1:B:246:TYR:CE2	1:B:275:PRO:HB2	2.47	0.50
1:A:332:THR:OG1	1:D:328:ARG:HG2	2.12	0.49
1:C:23:TRP:HA	1:C:26:LYS:HG3	1.95	0.49
1:A:97:ASP:HB3	1:A:115:VAL:HG13	1.94	0.49
1:A:85:ASN:HB2	1:A:86:PRO:HD2	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:142:LYS:HA	1:D:239:HIS:HB2	1.94	0.49
1:D:177:PRO:HD2	4:D:349:SO4:O3	2.12	0.49
1:A:274:ASN:HD22	1:A:277[B]:GLN:HG2	1.74	0.49
1:A:85:ASN:HB2	1:A:86:PRO:CD	2.42	0.49
1:C:38:GLU:CG	1:C:65:MET:CE	2.91	0.49
1:B:75:CYS:O	1:B:80:ALA:HB2	2.12	0.48
1:D:225:GLU:OE2	1:D:225:GLU:N	2.44	0.48
1:A:267:LEU:HD13	1:A:275:PRO:HB3	1.96	0.48
1:D:116:LEU:HD12	1:D:116:LEU:C	2.33	0.48
1:B:113:MET:HE1	1:B:303:MET:CE	2.41	0.48
1:C:113:MET:CE	1:C:303:MET:HE2	2.43	0.48
1:B:80:ALA:HB1	1:B:89:LEU:HD21	1.94	0.48
1:D:172:VAL:HG11	1:D:203:ALA:CB	2.44	0.48
1:A:279:TYR:HA	1:A:283:GLU:OE2	2.13	0.48
1:B:240:PHE:HB3	1:B:281:CYS:HB3	1.95	0.47
1:A:337:GLU:HA	1:A:338:SER:HA	1.46	0.47
1:D:23:TRP:HA	1:D:26:LYS:HD2	1.96	0.47
1:C:336:LYS:CB	7:C:781:HOH:O	2.62	0.47
1:D:115:VAL:HG22	1:D:116:LEU:N	2.29	0.47
1:B:50:ARG:HA	1:B:76:THR:HG23	1.97	0.47
1:C:10:ILE:CG1	1:C:319:VAL:HG21	2.44	0.47
1:C:128:PRO:HB2	1:C:130:PHE:CD2	2.50	0.47
1:A:273:LYS:CG	1:A:274:ASN:H	2.23	0.47
1:B:232:ALA:HB2	1:B:292:PHE:CD2	2.49	0.47
1:C:273:LYS:O	1:C:275:PRO:HD3	2.15	0.47
1:A:142:LYS:HA	1:A:239:HIS:HB2	1.96	0.46
1:B:44:MET:HE3	1:B:116:LEU:HD22	1.97	0.46
1:A:85:ASN:OD1	1:A:88:GLU:HG3	2.14	0.46
1:C:142:LYS:HE2	7:C:698:HOH:O	2.14	0.46
1:D:253:THR:HG22	1:D:258:GLU:OE1	2.14	0.46
1:A:55:ILE:HB	1:A:95:ALA:CB	2.46	0.46
1:B:277[B]:GLN:NE2	1:B:277[B]:GLN:CA	2.76	0.46
1:B:225:GLU:OE2	1:B:225:GLU:N	2.48	0.46
1:D:20:SER:HB2	1:D:36:ALA:HB2	1.98	0.45
1:C:113:MET:HE1	1:C:303:MET:HE2	1.98	0.45
1:B:78:GLU:HA	1:B:79:ASP:HA	1.51	0.45
1:A:34[B]:GLN:HG3	1:A:35:VAL:N	2.32	0.45
1:B:128:PRO:HB3	1:B:270:MET:CE	2.47	0.45
1:A:100:GLU:OE2	1:A:199:GLY:HA2	2.17	0.45
1:D:232:ALA:HB2	1:D:292:PHE:CD2	2.51	0.45
1:C:137:ALA:HB1	1:C:138:PRO:CD	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:287:GLY:O	5:A:355:CL:CL	2.72	0.45
1:D:248:PRO:HA	1:D:251:VAL:O	2.16	0.45
1:B:2:ASP:OD1	1:B:3:SER:N	2.49	0.45
1:B:138:PRO:HD3	1:B:215:HIS:O	2.17	0.45
1:B:270:MET:HB3	1:B:270:MET:HE2	1.55	0.45
5:A:363:CL:CL	7:A:678:HOH:O	2.58	0.44
1:B:177:PRO:O	5:B:353:CL:CL	2.72	0.44
1:C:116:LEU:C	1:C:116:LEU:HD12	2.37	0.44
1:D:10:ILE:HG12	1:D:319:VAL:HG21	1.99	0.44
1:A:274:ASN:O	1:A:277[B]:GLN:HB2	2.17	0.44
1:C:205:ILE:CG2	1:C:302:LEU:HD11	2.47	0.44
1:A:285:ALA:O	5:A:355:CL:CL	2.73	0.44
1:D:246:TYR:CE2	1:D:275:PRO:HB2	2.53	0.44
1:C:25:GLY:HA3	1:D:193:VAL:O	2.18	0.44
1:D:24:MET:HG3	1:D:108:GLY:HA2	1.99	0.44
1:A:101:GLY:CA	5:A:361:CL:CL	3.03	0.44
1:C:128:PRO:HB2	1:C:130:PHE:HD2	1.81	0.44
1:A:223:ALA:N	1:A:224:PRO:CD	2.80	0.44
1:B:10:ILE:HG12	1:B:319:VAL:HG21	2.00	0.43
1:B:77:ARG:N	1:B:77:ARG:HD3	2.32	0.43
1:C:164:ARG:NH2	3:C:348:AMP:O1P	2.47	0.43
1:A:37:VAL:HG22	1:A:98:PRO:HG2	2.00	0.43
1:B:50:ARG:NH1	1:B:76:THR:HA	2.34	0.43
1:A:139:PRO:HB2	1:D:340:LYS:O	2.18	0.43
1:D:223:ALA:N	1:D:224:PRO:CD	2.82	0.43
1:A:286:CSO:SG	7:A:457:HOH:O	2.62	0.43
1:A:255:LEU:HA	1:A:256:ILE:HA	1.63	0.43
1:C:256:ILE:HA	1:C:257:GLY:HA2	1.55	0.43
1:D:266:ARG:NH1	1:D:266:ARG:HG2	2.34	0.43
1:B:100:GLU:HG3	1:B:100:GLU:O	2.16	0.43
1:D:264:LEU:HA	1:D:264:LEU:HD12	1.87	0.42
1:A:116:LEU:C	1:A:116:LEU:HD12	2.39	0.42
1:C:67:TYR:CE1	1:C:70:GLU:HB2	2.55	0.42
1:A:164:ARG:HD2	1:A:168:GLU:O	2.20	0.42
1:C:113:MET:HE1	1:C:303:MET:CE	2.49	0.42
1:D:113:MET:SD	1:D:303:MET:HE1	2.59	0.42
1:D:172:VAL:HG11	1:D:203:ALA:HB1	2.00	0.42
1:C:282:GLU:O	1:C:286:CSO:HA	2.19	0.42
1:B:24:MET:CG	1:B:108:GLY:HA2	2.49	0.42
1:B:24:MET:HG3	1:B:108:GLY:HA2	2.01	0.42
1:B:223:ALA:HB3	1:B:224:PRO:HD3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:336:LYS:N	1:B:337:GLU:HA	2.35	0.42
1:D:248:PRO:O	1:D:252:LYS:HG2	2.19	0.42
1:D:223:ALA:HB3	1:D:224:PRO:HD3	2.01	0.41
1:C:3:SER:HA	1:C:323:GLN:OE1	2.19	0.41
1:A:267:LEU:O	1:A:272:ILE:HB	2.19	0.41
1:D:337:GLU:OE2	5:D:353:CL:CL	2.75	0.41
1:C:119:SER:OG	1:C:287:GLY:HA3	2.19	0.41
1:D:131:TYR:CE2	1:D:251:VAL:HG22	2.55	0.41
1:D:279:TYR:HA	1:D:283:GLU:OE2	2.20	0.41
1:A:267:LEU:HD23	1:A:270:MET:CE	2.50	0.41
1:B:74:ILE:O	1:B:77:ARG:HG2	2.20	0.41
1:A:114:ALA:O	1:A:294:ALA:HA	2.19	0.41
1:C:176:ARG:HB3	4:C:351:SO4:O2	2.21	0.41
1:C:225:GLU:OE2	1:C:225:GLU:N	2.53	0.41
1:B:267:LEU:HD13	1:B:275:PRO:HB3	2.01	0.41
1:C:102:THR:HG23	4:C:352:SO4:O4	2.21	0.41
1:C:138:PRO:HD3	1:C:215:HIS:O	2.21	0.41
1:C:252:LYS:HG2	1:C:254:GLY:H	1.86	0.41
1:B:49:MET:C	1:B:75:CYS:HB2	2.41	0.41
1:A:222:ALA:HB3	1:A:225:GLU:OE2	2.20	0.41
1:C:223:ALA:HB3	1:C:224:PRO:HD3	2.01	0.41
1:B:274[A]:ASN:HA	1:B:275:PRO:HD2	1.88	0.40
1:A:97:ASP:HB3	1:A:115:VAL:CG1	2.50	0.40
1:A:128:PRO:HB3	1:A:270:MET:SD	2.60	0.40
1:B:277[B]:GLN:NE2	1:B:278:VAL:N	2.65	0.40
1:B:223:ALA:N	1:B:224:PRO:CD	2.84	0.40
1:A:50:ARG:HA	1:A:76:THR:HG23	2.03	0.40
1:A:152:THR:O	1:A:156:LYS:HG3	2.22	0.40
1:A:119:SER:OG	1:A:287:GLY:HA3	2.22	0.40
1:B:311:GLY:HA3	1:B:339:PRO:HB3	2.03	0.40
1:B:336:LYS:HA	1:B:337:GLU:CB	2.52	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	348/379 (92%)	340 (98%)	7 (2%)	1 (0%)	46	57
1	B	346/379 (91%)	336 (97%)	10 (3%)	0	100	100
1	C	346/379 (91%)	340 (98%)	6 (2%)	0	100	100
1	D	345/379 (91%)	338 (98%)	7 (2%)	0	100	100
All	All	1385/1516 (91%)	1354 (98%)	30 (2%)	1 (0%)	56	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	273	LYS

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	272/301 (90%)	269 (99%)	3 (1%)	80	90
1	B	270/301 (90%)	268 (99%)	2 (1%)	88	95
1	C	266/301 (88%)	260 (98%)	6 (2%)	58	75
1	D	271/301 (90%)	267 (98%)	4 (2%)	72	85
All	All	1079/1204 (90%)	1064 (99%)	15 (1%)	74	86

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	SER
1	A	273	LYS
1	A	279	TYR
1	B	100	GLU
1	B	265	GLU
1	C	0	SER
1	C	34	GLN

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Mol	Chain	Res	Type
1	C	100	GLU
1	C	103	ASN
1	C	113	MET
1	C	338	SER
1	D	100	GLU
1	D	234	ARG
1	D	258	GLU
1	D	269	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	184	GLN
1	C	34	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](#)

4 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	CSO	A	286	1	3,6,7	0.52	0	1,6,8	1.92	0
1	CSO	B	286	1	3,6,7	0.53	0	1,6,8	1.98	0
1	CSO	C	286	1	3,6,7	0.52	0	1,6,8	1.98	0
1	CSO	D	286	1	3,6,7	0.52	0	1,6,8	1.88	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
1	CSO	A	286	1	-	0/1/5/7	0/0/0/0
1	CSO	B	286	1	-	0/1/5/7	0/0/0/0
1	CSO	C	286	1	-	0/1/5/7	0/0/0/0
1	CSO	D	286	1	-	0/1/5/7	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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	286	CSO	1	0
1	C	286	CSO	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 53 ligands modelled in this entry, 31 are monoatomic - leaving 22 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	AMP	A	348	-	20,25,25	0.93	1 (5%)	22,38,38	1.84	3 (13%)
4	SO4	A	349	-	4,4,4	0.24	0	6,6,6	0.07	0
4	SO4	A	350	-	4,4,4	0.20	0	6,6,6	0.09	0
4	SO4	A	351	-	4,4,4	0.22	0	6,6,6	0.09	0
4	SO4	A	352	-	4,4,4	0.20	0	6,6,6	0.08	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	A	353	-	4,4,4	0.21	0	6,6,6	0.10	0
4	SO4	A	354	-	4,4,4	0.21	0	6,6,6	0.09	0
6	GOL	A	365	-	5,5,5	0.33	0	5,5,5	0.71	0
3	AMP	B	348	-	20,25,25	0.89	1 (5%)	22,38,38	1.94	3 (13%)
4	SO4	B	349	-	4,4,4	0.23	0	6,6,6	0.07	0
4	SO4	B	350	-	4,4,4	0.23	0	6,6,6	0.07	0
3	AMP	C	348	-	20,25,25	0.88	1 (5%)	22,38,38	1.94	4 (18%)
3	AMP	C	349	-	20,25,25	0.91	1 (5%)	22,38,38	1.85	3 (13%)
4	SO4	C	350	-	4,4,4	0.26	0	6,6,6	0.14	0
4	SO4	C	351	-	4,4,4	0.21	0	6,6,6	0.11	0
4	SO4	C	352	-	4,4,4	0.23	0	6,6,6	0.08	0
4	SO4	C	353	-	4,4,4	0.21	0	6,6,6	0.10	0
4	SO4	D	348	-	4,4,4	0.24	0	6,6,6	0.24	0
4	SO4	D	349	-	4,4,4	0.21	0	6,6,6	0.06	0
4	SO4	D	350	-	4,4,4	0.22	0	6,6,6	0.08	0
4	SO4	D	351	-	4,4,4	0.22	0	6,6,6	0.09	0
4	SO4	D	352	-	4,4,4	0.20	0	6,6,6	0.09	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	AMP	A	348	-	-	0/6/26/26	0/3/3/3
4	SO4	A	349	-	-	0/0/0/0	0/0/0/0
4	SO4	A	350	-	-	0/0/0/0	0/0/0/0
4	SO4	A	351	-	-	0/0/0/0	0/0/0/0
4	SO4	A	352	-	-	0/0/0/0	0/0/0/0
4	SO4	A	353	-	-	0/0/0/0	0/0/0/0
4	SO4	A	354	-	-	0/0/0/0	0/0/0/0
6	GOL	A	365	-	-	0/4/4/4	0/0/0/0
3	AMP	B	348	-	-	0/6/26/26	0/3/3/3
4	SO4	B	349	-	-	0/0/0/0	0/0/0/0
4	SO4	B	350	-	-	0/0/0/0	0/0/0/0
3	AMP	C	348	-	-	0/6/26/26	0/3/3/3
3	AMP	C	349	-	-	0/6/26/26	0/3/3/3
4	SO4	C	350	-	-	0/0/0/0	0/0/0/0
4	SO4	C	351	-	-	0/0/0/0	0/0/0/0
4	SO4	C	352	-	-	0/0/0/0	0/0/0/0
4	SO4	C	353	-	-	0/0/0/0	0/0/0/0
4	SO4	D	348	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SO4	D	349	-	-	0/0/0/0	0/0/0/0
4	SO4	D	350	-	-	0/0/0/0	0/0/0/0
4	SO4	D	351	-	-	0/0/0/0	0/0/0/0
4	SO4	D	352	-	-	0/0/0/0	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	348	AMP	C5-C4	2.91	1.47	1.40
3	C	349	AMP	C5-C4	2.91	1.47	1.40
3	B	348	AMP	C5-C4	2.92	1.47	1.40
3	A	348	AMP	C5-C4	3.06	1.47	1.40

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	348	AMP	N3-C2-N1	-7.21	123.37	128.89
3	C	348	AMP	N3-C2-N1	-7.13	123.43	128.89
3	C	349	AMP	N3-C2-N1	-6.76	123.72	128.89
3	A	348	AMP	N3-C2-N1	-6.61	123.83	128.89
3	C	348	AMP	C4-C5-N7	-3.15	106.58	109.48
3	A	348	AMP	C4-C5-N7	-3.12	106.61	109.48
3	C	349	AMP	C4-C5-N7	-2.92	106.79	109.48
3	B	348	AMP	C4-C5-N7	-2.85	106.85	109.48
3	C	348	AMP	O3P-P-O5'	-2.10	100.52	106.56
3	C	349	AMP	O2P-P-O1P	2.34	118.13	110.58
3	B	348	AMP	O2P-P-O1P	2.34	118.13	110.58
3	C	348	AMP	O2P-P-O1P	2.35	118.13	110.58
3	A	348	AMP	O2P-P-O1P	2.35	118.15	110.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	365	GOL	2	0
3	C	348	AMP	1	0
4	C	351	SO4	1	0
4	C	352	SO4	1	0
4	D	349	SO4	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	346/379 (91%)	0.12	25 (7%) 18 26	17, 24, 61, 86	0
1	B	344/379 (90%)	0.20	32 (9%) 11 16	14, 25, 65, 93	0
1	C	346/379 (91%)	0.09	27 (7%) 16 22	15, 24, 71, 106	0
1	D	345/379 (91%)	0.18	28 (8%) 15 21	16, 26, 70, 90	0
All	All	1381/1516 (91%)	0.15	112 (8%) 15 21	14, 25, 67, 106	0

All (112) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	79	ASP	9.2
1	C	256	ILE	9.1
1	B	78	GLU	8.0
1	B	84	CYS	7.9
1	A	254	GLY	7.8
1	C	254	GLY	7.8
1	D	257	GLY	7.6
1	B	80	ALA	7.5
1	D	256	ILE	6.8
1	A	-1	GLY	6.7
1	C	255	LEU	6.6
1	B	256	ILE	6.6
1	D	254	GLY	6.5
1	D	255	LEU	6.3
1	B	82	SER	6.0
1	A	83	PHE	5.9
1	D	79	ASP	5.9
1	C	257	GLY	5.8
1	B	83	PHE	5.7
1	A	82	SER	5.7
1	B	254	GLY	5.3

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Mol	Chain	Res	Type	RSRZ
1	A	269	SER	5.3
1	C	269	SER	5.3
1	C	271	GLY	5.3
1	D	83	PHE	5.3
1	A	271	GLY	5.2
1	B	268	ALA	5.0
1	B	270	MET	4.9
1	D	82	SER	4.8
1	C	0	SER	4.6
1	D	86	PRO	4.6
1	B	257	GLY	4.5
1	A	79	ASP	4.4
1	D	78	GLU	4.4
1	A	256	ILE	4.3
1	C	268	ALA	4.3
1	D	269	SER	4.3
1	D	268	ALA	4.2
1	A	255	LEU	4.2
1	B	271	GLY	4.1
1	C	-1	GLY	4.1
1	C	261	GLU	4.0
1	D	85	ASN	4.0
1	B	75	CYS	4.0
1	D	80	ALA	3.9
1	C	253	THR	3.9
1	B	76	THR	3.9
1	A	257	GLY	3.9
1	A	268	ALA	3.7
1	A	0	SER	3.7
1	C	249	GLU	3.7
1	D	62	ASP	3.7
1	A	80	ALA	3.7
1	D	270	MET	3.7
1	D	273	LYS	3.7
1	D	77	ARG	3.5
1	A	78	GLU	3.5
1	D	0	SER	3.5
1	B	261	GLU	3.4
1	B	266	ARG	3.4
1	B	252	LYS	3.4
1	B	255	LEU	3.4
1	B	77	ARG	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	77	ARG	3.4
1	A	86	PRO	3.4
1	A	76	THR	3.3
1	B	262	GLY	3.3
1	D	261	GLU	3.3
1	D	61	ASP	3.3
1	B	273	LYS	3.3
1	A	272	ILE	3.3
1	B	1	VAL	3.3
1	B	258	GLU	3.2
1	C	266	ARG	3.2
1	C	270	MET	3.2
1	B	274[A]	ASN	3.1
1	A	267	LEU	3.1
1	D	266	ARG	3.0
1	C	258	GLU	3.0
1	B	259	SER	2.9
1	A	274	ASN	2.8
1	C	60	ARG	2.8
1	A	270	MET	2.8
1	C	336	LYS	2.8
1	A	337	GLU	2.8
1	C	267	LEU	2.7
1	B	86	PRO	2.7
1	D	60	ARG	2.7
1	C	259	SER	2.6
1	B	85	ASN	2.6
1	A	248	PRO	2.6
1	D	252	LYS	2.6
1	B	265	GLU	2.6
1	B	269	SER	2.6
1	A	261	GLU	2.6
1	C	265	GLU	2.6
1	D	271	GLY	2.6
1	B	272	ILE	2.5
1	D	337	GLU	2.5
1	C	62	ASP	2.5
1	D	274	ASN	2.4
1	C	252	LYS	2.4
1	C	273	LYS	2.3
1	B	253	THR	2.3
1	C	1	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	82	SER	2.3
1	C	272	ILE	2.3
1	B	81	LYS	2.2
1	D	265	GLU	2.2
1	A	273	LYS	2.2
1	D	250	VAL	2.1
1	C	61	ASP	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	CSO	D	286	7/8	0.94	0.12	-	29,30,34,36	0
1	CSO	C	286	7/8	0.96	0.19	-	26,27,32,57	0
1	CSO	B	286	7/8	0.97	0.08	-	25,25,30,31	0
1	CSO	A	286	7/8	0.96	0.13	-	25,26,29,54	0

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(Å ²)	Q<0.9
2	MG	A	347	1/1	0.97	0.22	19.02	36,36,36,36	0
2	MG	B	347	1/1	0.97	0.23	8.91	36,36,36,36	0
4	SO4	D	350	5/5	0.96	0.23	7.19	37,44,52,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	SO4	C	352	5/5	0.98	0.25	6.70	44,45,54,60	0
6	GOL	A	365	6/6	0.83	0.21	5.23	33,40,50,52	0
4	SO4	A	350	5/5	0.97	0.28	2.98	44,50,54,57	0
2	MG	C	347	1/1	0.95	0.19	2.70	39,39,39,39	0
5	CL	A	355	1/1	0.95	0.23	2.58	45,45,45,45	0
5	CL	A	358	1/1	0.77	0.20	2.51	55,55,55,55	0
4	SO4	B	350	5/5	0.98	0.23	1.60	37,51,52,52	0
5	CL	A	363	1/1	0.78	0.14	0.90	50,50,50,50	0
4	SO4	D	349	5/5	0.97	0.24	0.69	37,42,46,50	0
2	MG	D	347	1/1	0.96	0.11	0.59	36,36,36,36	0
3	AMP	B	348	23/23	0.98	0.11	-0.36	15,17,18,19	0
3	AMP	C	349	23/23	0.98	0.10	-0.38	16,18,19,20	0
4	SO4	C	350	5/5	0.97	0.11	-0.40	37,39,43,47	0
3	AMP	A	348	23/23	0.98	0.10	-0.59	16,18,19,22	0
3	AMP	C	348	23/23	0.98	0.10	-0.70	15,16,17,18	0
5	CL	D	353	1/1	0.81	0.09	-0.83	53,53,53,53	0
4	SO4	D	348	5/5	0.98	0.09	-1.35	37,39,42,43	0
5	CL	C	356	1/1	0.99	0.06	-1.86	38,38,38,38	0
5	CL	A	361	1/1	0.36	0.20	-	55,55,55,55	0
5	CL	B	352	1/1	0.92	0.08	-	71,71,71,71	0
4	SO4	A	353	5/5	0.95	0.29	-	55,55,69,72	0
5	CL	C	354	1/1	0.87	0.23	-	59,59,59,59	0
5	CL	A	359	1/1	0.94	0.12	-	55,55,55,55	0
5	CL	C	355	1/1	0.71	0.16	-	58,58,58,58	0
4	SO4	D	352	5/5	0.92	0.31	-	47,61,71,78	0
5	CL	C	357	1/1	0.91	0.12	-	61,61,61,61	0
4	SO4	B	349	5/5	0.98	0.08	-	35,36,42,43	0
5	CL	B	355	1/1	0.81	0.59	-	115,115,115,115	0
4	SO4	D	351	5/5	0.96	0.30	-	42,52,54,60	0
2	MG	C	346	1/1	0.95	0.26	-	42,42,42,42	0
5	CL	A	356	1/1	0.78	0.25	-	61,61,61,61	0
4	SO4	A	349	5/5	0.99	0.14	-	34,37,39,43	0
2	MG	B	346	1/1	0.84	0.18	-	48,48,48,48	0
5	CL	C	359	1/1	0.73	0.14	-	57,57,57,57	0
5	CL	B	353	1/1	0.38	0.23	-	57,57,57,57	0
4	SO4	A	351	5/5	0.95	0.22	-	42,52,54,62	0
2	MG	A	346	1/1	0.89	0.34	-	52,52,52,52	0
5	CL	A	357	1/1	0.80	0.22	-	53,53,53,53	0
5	CL	A	364	1/1	0.82	0.09	-	57,57,57,57	0
5	CL	B	351	1/1	0.95	0.09	-	49,49,49,49	0
4	SO4	C	351	5/5	0.98	0.32	-	40,46,49,53	0
5	CL	A	362	1/1	0.89	0.33	-	66,66,66,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	SO4	C	353	5/5	0.94	0.31	-	52,55,63,67	0
2	MG	D	346	1/1	0.96	0.29	-	44,44,44,44	0
5	CL	D	354	1/1	0.74	0.13	-	74,74,74,74	0
4	SO4	A	352	5/5	0.96	0.25	-	56,57,63,67	0
5	CL	B	354	1/1	0.92	0.06	-	53,53,53,53	0
5	CL	C	358	1/1	0.75	0.18	-	67,67,67,67	0
4	SO4	A	354	5/5	0.93	0.30	-	54,58,69,72	0
5	CL	A	360	1/1	0.86	0.12	-	65,65,65,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.