Full wwPDB X-ray Structure Validation Report (i)

Feb 15, 2024 – 01:47 AM EST

PDB ID	:	3NIY
Title	:	Crystal structure of native xylanase 10B from Thermotoga petrophila RKU-1
Authors	:	Santos, C.R.; Meza, A.N.; Trindade, D.M.; Ruller, R.; Squina, F.M.; Prade, R.A.; Murakami, M.T.
Deposited on	:	2010-06-16
Resolution	:	1.58 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

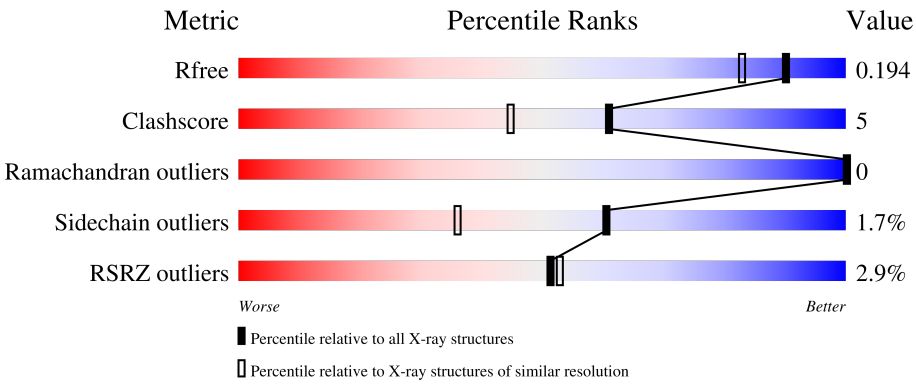
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.58 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5534 (1.60-1.56)
Clashscore	141614	5861 (1.60-1.56)
Ramachandran outliers	138981	5708 (1.60-1.56)
Sidechain outliers	138945	5703 (1.60-1.56)
RSRZ outliers	127900	5431 (1.60-1.56)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	341	3% 85% 9% • •
1	B	341	3% 81% 12% • 5%

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
3	ACT	A	347	-	X	-	-
3	ACT	B	343	-	-	X	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6113 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 Endo-1,4-beta-xylanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	326	2736	1767	453	506	10	0	4	0
1	B	325	2716	1757	448	501	10	0	1	0

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP A5IL00
A	2	GLY	-	expression tag	UNP A5IL00
A	3	HIS	-	expression tag	UNP A5IL00
A	4	HIS	-	expression tag	UNP A5IL00
A	5	HIS	-	expression tag	UNP A5IL00
A	6	HIS	-	expression tag	UNP A5IL00
A	7	HIS	-	expression tag	UNP A5IL00
A	8	HIS	-	expression tag	UNP A5IL00
A	9	GLU	-	expression tag	UNP A5IL00
A	10	ASN	-	expression tag	UNP A5IL00
A	11	LEU	-	expression tag	UNP A5IL00
A	12	TYR	-	expression tag	UNP A5IL00
A	13	PHE	-	expression tag	UNP A5IL00
A	14	GLN	-	expression tag	UNP A5IL00
A	15	GLY	-	expression tag	UNP A5IL00
A	16	HIS	-	expression tag	UNP A5IL00
A	17	MET	-	expression tag	UNP A5IL00
B	1	MET	-	expression tag	UNP A5IL00
B	2	GLY	-	expression tag	UNP A5IL00
B	3	HIS	-	expression tag	UNP A5IL00
B	4	HIS	-	expression tag	UNP A5IL00
B	5	HIS	-	expression tag	UNP A5IL00
B	6	HIS	-	expression tag	UNP A5IL00
B	7	HIS	-	expression tag	UNP A5IL00
B	8	HIS	-	expression tag	UNP A5IL00

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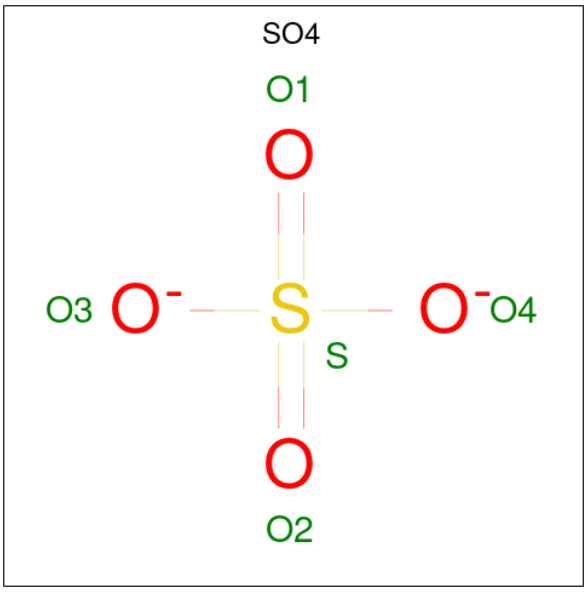
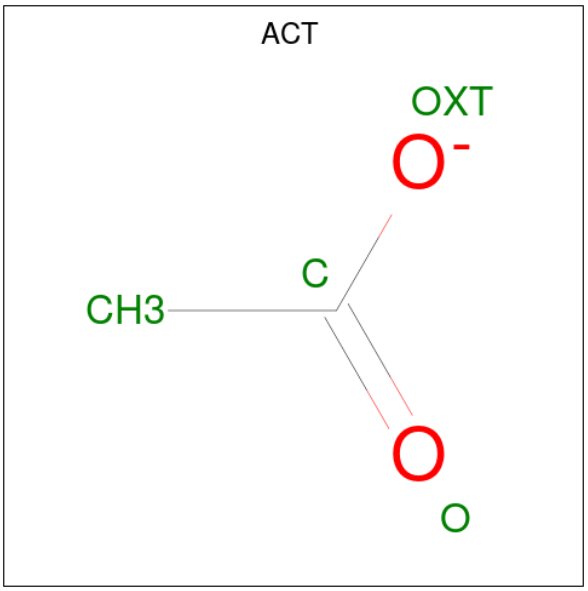
Chain	Residue	Modelled	Actual	Comment	Reference
B	9	GLU	-	expression tag	UNP A5IL00
B	10	ASN	-	expression tag	UNP A5IL00
B	11	LEU	-	expression tag	UNP A5IL00
B	12	TYR	-	expression tag	UNP A5IL00
B	13	PHE	-	expression tag	UNP A5IL00
B	14	GLN	-	expression tag	UNP A5IL00
B	15	GLY	-	expression tag	UNP A5IL00
B	16	HIS	-	expression tag	UNP A5IL00
B	17	MET	-	expression tag	UNP A5IL00

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 4 is water.

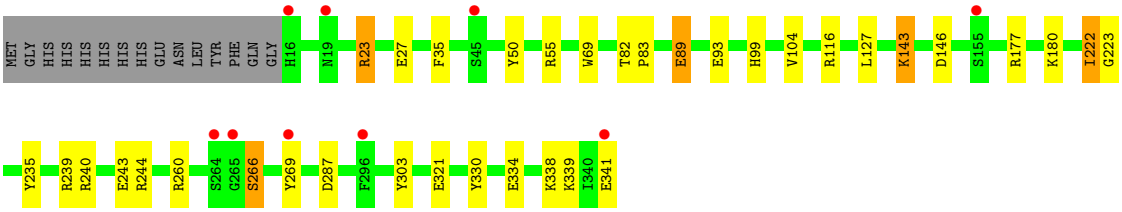
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	316	316	316	0	0
			Total	O		
4	B	300	300	300	0	0

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.

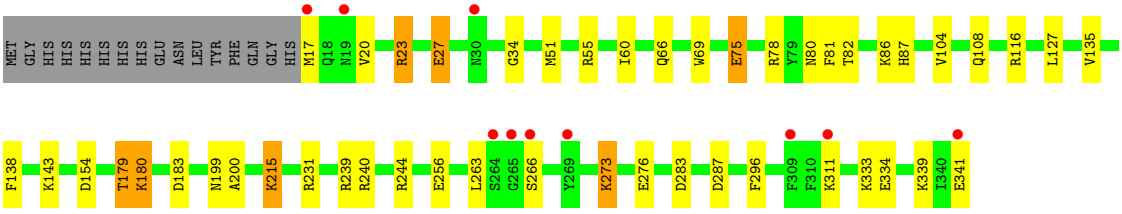
- Molecule 1: Endo-1,4-beta-xylanase

Chain A:



- Molecule 1: Endo-1,4-beta-xylanase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	58.54Å 58.55Å 61.58Å 84.54° 70.82° 68.87°	Depositor
Resolution (Å)	31.24 – 1.58 31.24 – 1.58	Depositor EDS
% Data completeness (in resolution range)	92.5 (31.24-1.58) 92.5 (31.24-1.58)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.33 (at 1.58Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.158 , 0.196 0.156 , 0.194	Depositor DCC
R_{free} test set	4584 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	14.6	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 46.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	6113	wwPDB-VP
Average B, all atoms (Å^2)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.77% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: ACT, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\|>5$	RMSZ	$\#\|Z\|>5$
1	A	1.41	7/2831 (0.2%)	1.27	11/3831 (0.3%)
1	B	1.46	14/2796 (0.5%)	1.28	14/3783 (0.4%)
All	All	1.43	21/5627 (0.4%)	1.27	25/7614 (0.3%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	75	GLU	CG-CD	-10.64	1.35	1.51
1	B	75	GLU	CB-CG	-9.93	1.33	1.52
1	B	215	LYS	CB-CG	-9.07	1.28	1.52
1	B	334	GLU	CG-CD	8.84	1.65	1.51
1	A	69	TRP	CB-CG	-8.45	1.35	1.50
1	B	69	TRP	CB-CG	-7.84	1.36	1.50
1	B	334	GLU	CD-OE2	7.73	1.34	1.25
1	A	50	TYR	CD1-CE1	7.21	1.50	1.39
1	B	27	GLU	CG-CD	6.70	1.61	1.51
1	A	239	ARG	CG-CD	6.67	1.68	1.51
1	B	240	ARG	CG-CD	6.38	1.68	1.51
1	B	86	LYS	CD-CE	6.32	1.67	1.51
1	B	215	LYS	CA-CB	-5.71	1.41	1.53
1	A	235	TYR	CG-CD2	5.70	1.46	1.39
1	B	256	GLU	CB-CG	5.67	1.62	1.52
1	B	55	ARG	CG-CD	5.61	1.66	1.51
1	B	180	LYS	CE-NZ	5.60	1.63	1.49
1	A	55	ARG	CG-CD	5.38	1.65	1.51
1	A	243	GLU	CD-OE2	5.27	1.31	1.25
1	A	89	GLU	CG-CD	5.14	1.59	1.51
1	B	296	PHE	CD1-CE1	5.10	1.49	1.39

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	51	MET	CG-SD-CE	-8.70	86.29	100.20
1	B	116	ARG	NE-CZ-NH2	-8.24	116.18	120.30
1	A	116	ARG	NE-CZ-NH1	-7.98	116.31	120.30
1	B	244	ARG	NE-CZ-NH2	-7.96	116.32	120.30
1	B	23	ARG	NE-CZ-NH2	-7.27	116.67	120.30
1	A	177	ARG	NE-CZ-NH2	-7.11	116.74	120.30
1	A	244	ARG	NE-CZ-NH2	-6.97	116.81	120.30
1	A	50	TYR	CD1-CE1-CZ	-6.93	113.56	119.80
1	B	334	GLU	CG-CD-OE2	6.84	131.97	118.30
1	B	138	PHE	CB-CG-CD2	-6.63	116.16	120.80
1	A	239	ARG	CG-CD-NE	6.57	125.60	111.80
1	B	215	LYS	CB-CG-CD	-6.29	95.25	111.60
1	A	35	PHE	CB-CG-CD2	-6.08	116.55	120.80
1	A	303	TYR	CB-CG-CD2	-5.78	117.53	121.00
1	A	239	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	A	116	ARG	NE-CZ-NH2	5.69	123.14	120.30
1	B	283	ASP	CB-CG-OD2	-5.64	113.22	118.30
1	A	260	ARG	NE-CZ-NH1	5.62	123.11	120.30
1	B	244	ARG	NH1-CZ-NH2	5.48	125.43	119.40
1	B	81	PHE	CB-CG-CD2	-5.46	116.97	120.80
1	B	296	PHE	CG-CD1-CE1	-5.30	114.97	120.80
1	B	55	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	A	23	ARG	NE-CZ-NH2	-5.21	117.69	120.30
1	B	179	THR	CA-CB-CG2	-5.08	105.29	112.40
1	B	116	ARG	NE-CZ-NH1	5.01	122.81	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2736	0	2647	18	0
1	B	2716	0	2636	31	0
2	A	20	0	0	0	1
2	B	5	0	0	0	0
3	A	12	0	9	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	8	0	6	4	0
4	A	316	0	0	6	1
4	B	300	0	0	8	2
All	All	6113	0	5298	49	3

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 5.

All (49) 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:B:180:LYS:HD2	4:B:555:HOH:O	1.51	1.09
1:B:108:GLN:NE2	4:B:437:HOH:O	1.89	1.05
1:B:154:ASP:O	4:B:459:HOH:O	1.83	0.96
1:A:341:GLU:OXT	1:A:341:GLU:HG3	1.74	0.87
1:B:183:ASP:OD2	4:B:546:HOH:O	2.06	0.74
1:A:321:GLU:H	1:A:321:GLU:CD	1.91	0.73
4:A:621:HOH:O	1:B:78:ARG:CG	2.37	0.71
1:A:330:TYR:O	1:A:334:GLU:HG3	1.97	0.65
1:A:180:LYS:HD2	4:A:610:HOH:O	1.98	0.63
1:A:82:THR:HB	1:A:83:PRO:HD3	1.81	0.62
1:B:27:GLU:OE2	4:B:643:HOH:O	2.15	0.62
1:B:200:ALA:H	3:B:343:ACT:CH3	2.14	0.60
1:B:273:LYS:CE	1:B:276:GLU:OE1	2.49	0.60
1:B:231:ARG:HH11	1:B:231:ARG:HG2	1.67	0.59
1:B:273:LYS:HE3	1:B:276:GLU:OE1	2.03	0.59
1:A:27:GLU:OE1	4:A:419:HOH:O	2.17	0.57
1:B:66:GLN:OE1	1:B:87:HIS:HE1	1.91	0.54
1:B:273:LYS:HE3	1:B:273:LYS:HA	1.90	0.54
1:A:341:GLU:OXT	1:A:341:GLU:CG	2.49	0.53
1:A:338:LYS:HA	1:A:341:GLU:HG2	1.91	0.52
1:B:231:ARG:HG2	1:B:231:ARG:NH1	2.24	0.50
1:B:143:LYS:HG3	4:B:408:HOH:O	2.12	0.49
1:B:34:GLY:HA3	1:B:60:ILE:HG23	1.95	0.49
4:A:621:HOH:O	1:B:78:ARG:HG3	2.10	0.49
1:A:143:LYS:HG3	4:A:402:HOH:O	2.13	0.48
1:A:321:GLU:CD	1:A:321:GLU:N	2.65	0.48
1:A:89:GLU:OE1	1:A:93:GLU:OE1	2.32	0.47
1:B:200:ALA:H	3:B:343:ACT:H1	1.79	0.47
4:A:621:HOH:O	1:B:78:ARG:HD2	2.14	0.47
1:A:287:ASP:OD2	1:A:339:LYS:NZ	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:VAL:O	1:B:333:LYS:CE	2.64	0.46
1:A:23:ARG:O	1:A:27:GLU:HG3	2.16	0.45
1:A:104:VAL:HG22	1:A:127:LEU:HD11	1.99	0.45
1:B:23:ARG:O	1:B:27:GLU:HG3	2.17	0.45
1:B:239:ARG:HD3	4:B:615:HOH:O	2.18	0.43
1:B:200:ALA:H	3:B:343:ACT:H2	1.83	0.43
1:A:240:ARG:HH11	1:A:240:ARG:HD3	1.64	0.43
1:A:266:SER:HB2	1:A:269:TYR:HB3	2.01	0.43
1:B:273:LYS:HE2	1:B:276:GLU:OE1	2.19	0.42
1:A:99:HIS:NE2	1:A:146:ASP:OD1	2.52	0.42
1:B:199:ASN:HB2	3:B:343:ACT:H1	2.01	0.42
1:B:20:VAL:O	1:B:333:LYS:HE2	2.20	0.42
1:B:108:GLN:O	4:B:436:HOH:O	2.22	0.41
1:B:104:VAL:HG22	1:B:127:LEU:HD11	2.02	0.41
1:B:263:LEU:HD23	1:B:263:LEU:HA	1.77	0.41
1:A:222:ILE:HG13	1:A:223:GLY:N	2.36	0.41
1:B:80:ASN:ND2	1:B:82:THR:H	2.19	0.41
1:B:287:ASP:OD2	1:B:339:LYS:NZ	2.54	0.41
1:B:135:VAL:HG21	1:B:179:THR:HG23	2.03	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:550:HOH:O	4:B:616:HOH:O[1_455]	2.11	0.09
2:A:345:SO4:O2	4:B:629:HOH:O[1_544]	2.12	0.08
4:A:538:HOH:O	4:A:650:HOH:O[1_655]	2.18	0.02

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	328/341 (96%)	320 (98%)	8 (2%)	0	100	100
1	B	324/341 (95%)	318 (98%)	6 (2%)	0	100	100
All	All	652/682 (96%)	638 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	295/304 (97%)	292 (99%)	3 (1%)	76	59
1	B	291/304 (96%)	284 (98%)	7 (2%)	49	22
All	All	586/608 (96%)	576 (98%)	10 (2%)	60	36

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	143	LYS
1	A	222	ILE
1	A	266	SER
1	B	17	MET
1	B	75	GLU
1	B	215	LYS
1	B	266	SER
1	B	273	LYS
1	B	311	LYS
1	B	341	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	251	GLN
1	B	80	ASN
1	B	87	HIS

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Mol	Chain	Res	Type
1	B	95	ASN
1	B	251	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

10 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	$\#\|Z\| > 2$
3	ACT	B	344	-	3,3,3	1.37	0	3,3,3	1.82	1 (33%)
2	SO4	A	345	-	4,4,4	0.78	0	6,6,6	0.56	0
3	ACT	A	346	-	3,3,3	1.27	0	3,3,3	2.09	1 (33%)
2	SO4	A	343	-	4,4,4	0.51	0	6,6,6	0.78	0
3	ACT	B	343	-	3,3,3	1.36	0	3,3,3	2.23	2 (66%)
2	SO4	A	342	-	4,4,4	0.83	0	6,6,6	0.89	0
3	ACT	A	347	-	3,3,3	1.77	1 (33%)	3,3,3	2.35	2 (66%)
2	SO4	B	342	-	4,4,4	0.24	0	6,6,6	0.55	0
2	SO4	A	344	-	4,4,4	0.53	0	6,6,6	1.09	0
3	ACT	A	348	-	3,3,3	0.87	0	3,3,3	1.39	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	347	ACT	OXT-C	2.09	1.40	1.30

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	347	ACT	OXT-C-CH3	3.05	127.77	115.18
3	A	346	ACT	OXT-C-O	-3.00	111.00	122.05
3	B	343	ACT	OXT-C-CH3	2.76	126.58	115.18
3	B	343	ACT	OXT-C-O	-2.68	112.17	122.05
3	A	347	ACT	O-C-CH3	-2.64	112.04	122.33
3	B	344	ACT	OXT-C-O	-2.42	113.14	122.05

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	345	SO4	0	1
3	B	343	ACT	4	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	326/341 (95%)	-0.05	9 (2%)	53	54	8, 14, 24, 44	11 (3%)
1	B	325/341 (95%)	-0.07	10 (3%)	49	50	8, 14, 26, 47	7 (2%)
All	All	651/682 (95%)	-0.06	19 (2%)	51	53	8, 14, 26, 47	18 (2%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	265	GLY	4.6
1	A	265	GLY	4.3
1	B	264	SER	3.8
1	A	264	SER	3.4
1	A	341	GLU	3.2
1	A	16	HIS	3.0
1	B	30	ASN	2.9
1	B	341	GLU	2.8
1	B	17	MET	2.8
1	B	309	PHE	2.6
1	A	19	ASN	2.5
1	B	19	ASN	2.4
1	B	311	LYS	2.3
1	A	296	PHE	2.3
1	A	269	TYR	2.1
1	B	269	TYR	2.1
1	A	45	SER	2.1
1	B	266	SER	2.0
1	A	155	SER	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 monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	ACT	B	343	4/4	0.84	0.14	10,22,28,31	0
3	ACT	A	346	4/4	0.89	0.14	36,38,39,39	0
3	ACT	A	347	4/4	0.90	0.13	18,25,28,29	0
3	ACT	A	348	4/4	0.91	0.11	22,24,25,27	0
2	SO4	A	343	5/5	0.93	0.20	40,43,45,48	0
3	ACT	B	344	4/4	0.93	0.11	26,27,27,29	0
2	SO4	A	345	5/5	0.94	0.18	36,41,43,44	0
2	SO4	A	344	5/5	0.95	0.20	32,34,38,43	0
2	SO4	B	342	5/5	0.97	0.19	50,51,51,54	0
2	SO4	A	342	5/5	0.97	0.22	25,34,38,40	0

6.5 Other polymers (i)

There are no such residues in this entry.