



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:23 PM GMT

PDB ID : 3XIS  
Title : A METAL-MEDIATED HYDRIDE SHIFT MECHANISM FOR XYLOSE ISOMERASE BASED ON THE 1.6 ANGSTROMS STREPTOMYCES RUBIGINOSUS STRUCTURES WITH XYLITOL AND D-XYLOSE  
Authors : Whitlow, M.; Howard, A.J.  
Deposited on : 1991-03-25  
Resolution : 1.60 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

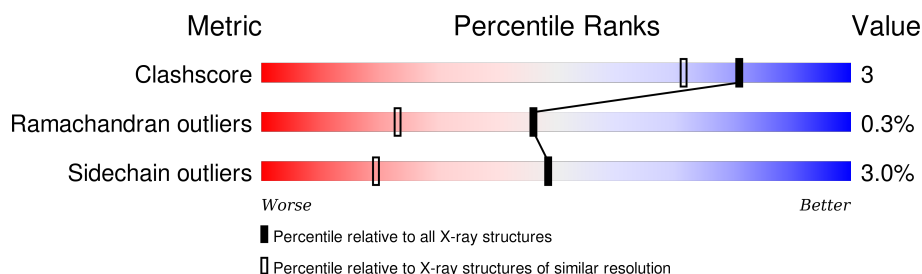
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.60 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	2732 (1.60-1.60)
Ramachandran outliers	100387	2654 (1.60-1.60)
Sidechain outliers	100360	2653 (1.60-1.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	387	

## 2 Entry composition [i](#)

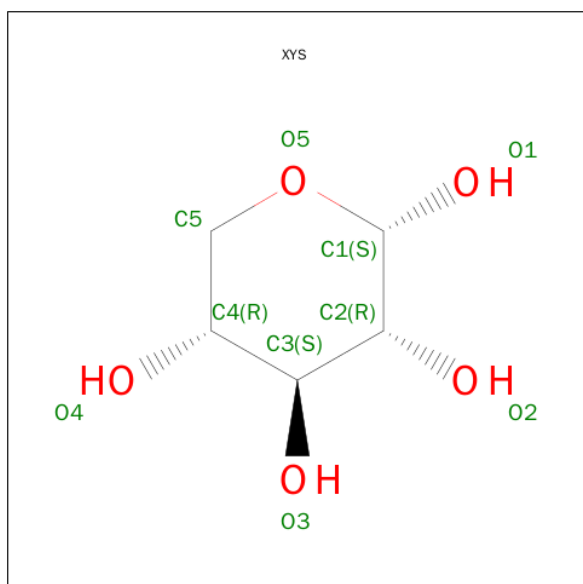
There are 4 unique types of molecules in this entry. The entry contains 3433 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 XYLOSE ISOMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	386	3049	1913	558	569	9	0	6	1

- Molecule 2 is SUGAR (2-MER) (three-letter code: XYS, XLS) (formula: C<sub>5</sub>H<sub>10</sub>O<sub>5</sub>, C<sub>5</sub>H<sub>10</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	20	10	10	0	1

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is water.

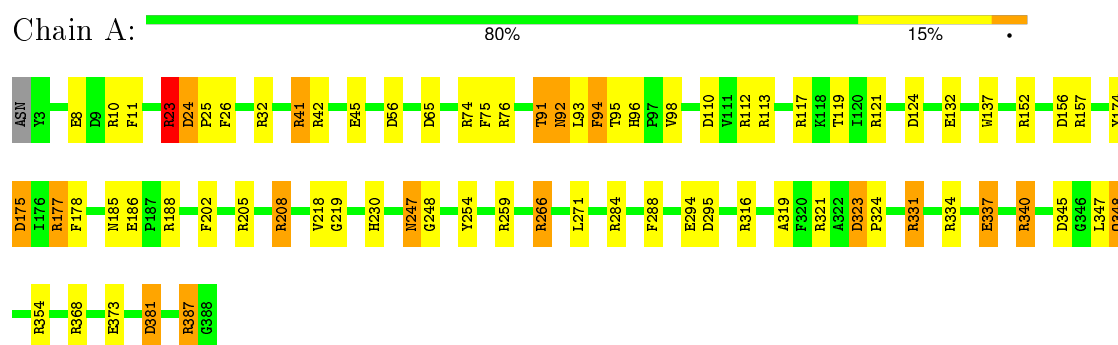
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	362	Total 362	O 362	0	0

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

#### • Molecule 1: XYLOSE ISOMERASE



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.64Å 99.97Å 103.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.60	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-1.60)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, $R_{free}$	0.151 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3433	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	17.0	wwPDB-VP

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MG, XYS, XLS

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	1.29	4/3158 (0.1%)	1.90	95/4270 (2.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	7

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	132[A]	GLU	CG-CD	-7.01	1.41	1.51
1	A	132[B]	GLU	CG-CD	-7.01	1.41	1.51
1	A	373	GLU	CD-OE1	-6.51	1.18	1.25
1	A	340	ARG	CD-NE	-5.85	1.36	1.46

All (95) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	331	ARG	NE-CZ-NH1	19.02	129.81	120.30
1	A	132[A]	GLU	CB-CG-CD	16.84	159.68	114.20
1	A	132[B]	GLU	CB-CG-CD	16.84	159.68	114.20
1	A	340	ARG	NE-CZ-NH1	16.39	128.50	120.30
1	A	188	ARG	NE-CZ-NH1	16.05	128.32	120.30
1	A	340	ARG	CD-NE-CZ	14.44	143.82	123.60
1	A	340	ARG	NE-CZ-NH2	-14.38	113.11	120.30
1	A	266	ARG	CD-NE-CZ	13.62	142.67	123.60
1	A	113	ARG	NE-CZ-NH1	12.37	126.49	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	188	ARG	NE-CZ-NH2	-11.48	114.56	120.30
1	A	152	ARG	CD-NE-CZ	11.06	139.09	123.60
1	A	266	ARG	NE-CZ-NH2	-10.97	114.81	120.30
1	A	368	ARG	NE-CZ-NH1	10.70	125.65	120.30
1	A	23	ARG	CD-NE-CZ	10.69	138.56	123.60
1	A	331	ARG	NE-CZ-NH2	-10.66	114.97	120.30
1	A	124	ASP	CB-CG-OD1	10.18	127.46	118.30
1	A	152	ARG	NE-CZ-NH1	9.29	124.95	120.30
1	A	177	ARG	NE-CZ-NH2	-9.22	115.69	120.30
1	A	56	ASP	CB-CG-OD1	9.21	126.59	118.30
1	A	74	ARG	CD-NE-CZ	9.07	136.29	123.60
1	A	266	ARG	NE-CZ-NH1	9.00	124.80	120.30
1	A	23	ARG	CG-CD-NE	8.98	130.66	111.80
1	A	76[A]	ARG	NE-CZ-NH2	-8.86	115.87	120.30
1	A	76[B]	ARG	NE-CZ-NH2	-8.86	115.87	120.30
1	A	74	ARG	NE-CZ-NH1	8.73	124.67	120.30
1	A	381	ASP	CB-CG-OD1	8.32	125.79	118.30
1	A	74	ARG	NE-CZ-NH2	-8.22	116.19	120.30
1	A	387[A]	ARG	NE-CZ-NH1	8.03	124.31	120.30
1	A	387[B]	ARG	NE-CZ-NH1	8.03	124.31	120.30
1	A	76[A]	ARG	NE-CZ-NH1	7.98	124.29	120.30
1	A	76[B]	ARG	NE-CZ-NH1	7.98	124.29	120.30
1	A	10	ARG	NE-CZ-NH1	7.94	124.27	120.30
1	A	157	ARG	NE-CZ-NH2	7.69	124.14	120.30
1	A	284	ARG	NE-CZ-NH1	7.68	124.14	120.30
1	A	42	ARG	NE-CZ-NH1	7.62	124.11	120.30
1	A	208	ARG	NE-CZ-NH2	-7.40	116.60	120.30
1	A	295	ASP	CB-CG-OD1	7.32	124.88	118.30
1	A	121	ARG	NE-CZ-NH2	-7.23	116.68	120.30
1	A	205[A]	ARG	NE-CZ-NH2	-7.21	116.70	120.30
1	A	205[B]	ARG	NE-CZ-NH2	-7.21	116.70	120.30
1	A	42	ARG	NE-CZ-NH2	-7.15	116.73	120.30
1	A	284	ARG	NE-CZ-NH2	-7.12	116.74	120.30
1	A	331	ARG	CD-NE-CZ	6.92	133.28	123.60
1	A	113	ARG	NE-CZ-NH2	-6.91	116.84	120.30
1	A	205[A]	ARG	NE-CZ-NH1	6.89	123.75	120.30
1	A	205[B]	ARG	NE-CZ-NH1	6.89	123.75	120.30
1	A	177	ARG	NE-CZ-NH1	6.87	123.74	120.30
1	A	368	ARG	CD-NE-CZ	6.80	133.12	123.60
1	A	41	ARG	NE-CZ-NH1	6.79	123.69	120.30
1	A	294	GLU	O-C-N	6.71	133.44	122.70
1	A	340	ARG	CG-CD-NE	6.59	125.63	111.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	45	GLU	CA-CB-CG	6.54	127.80	113.40
1	A	32	ARG	NE-CZ-NH1	6.42	123.51	120.30
1	A	259	ARG	NE-CZ-NH1	6.34	123.47	120.30
1	A	121	ARG	NE-CZ-NH1	6.31	123.45	120.30
1	A	354	ARG	CD-NE-CZ	6.18	132.25	123.60
1	A	174	TYR	CA-CB-CG	6.17	125.12	113.40
1	A	345	ASP	CB-CG-OD1	6.15	123.83	118.30
1	A	368	ARG	CA-CB-CG	6.04	126.70	113.40
1	A	23	ARG	NE-CZ-NH1	6.00	123.30	120.30
1	A	8	GLU	CA-CB-CG	5.99	126.58	113.40
1	A	23	ARG	NE-CZ-NH2	-5.80	117.40	120.30
1	A	323	ASP	CB-CG-OD2	5.80	123.52	118.30
1	A	156	ASP	CB-CG-OD2	-5.74	113.14	118.30
1	A	381	ASP	CB-CG-OD2	-5.71	113.16	118.30
1	A	205[A]	ARG	CD-NE-CZ	5.70	131.58	123.60
1	A	205[B]	ARG	CD-NE-CZ	5.70	131.58	123.60
1	A	94	PHE	CB-CG-CD1	-5.68	116.82	120.80
1	A	112	ARG	NE-CZ-NH2	-5.66	117.47	120.30
1	A	202	PHE	CB-CG-CD1	-5.63	116.86	120.80
1	A	10	ARG	CA-CB-CG	5.61	125.74	113.40
1	A	110	ASP	CB-CG-OD1	5.61	123.35	118.30
1	A	348	GLN	N-CA-CB	5.56	120.61	110.60
1	A	316	ARG	NE-CZ-NH2	-5.56	117.52	120.30
1	A	65	ASP	CB-CG-OD1	5.54	123.28	118.30
1	A	288	PHE	CB-CG-CD1	-5.48	116.97	120.80
1	A	75	PHE	CB-CG-CD1	-5.46	116.98	120.80
1	A	387[A]	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	A	387[B]	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	A	316	ARG	NE-CZ-NH1	5.44	123.02	120.30
1	A	124	ASP	CB-CG-OD2	-5.40	113.44	118.30
1	A	334	ARG	NE-CZ-NH1	5.39	122.99	120.30
1	A	337	GLU	CG-CD-OE2	-5.36	107.58	118.30
1	A	271	LEU	CB-CG-CD2	-5.33	101.95	111.00
1	A	24	ASP	CB-CA-C	5.31	121.03	110.40
1	A	254	TYR	CB-CG-CD2	-5.28	117.83	121.00
1	A	319	ALA	CB-CA-C	5.21	117.91	110.10
1	A	10	ARG	NE-CZ-NH2	-5.21	117.70	120.30
1	A	354	ARG	NE-CZ-NH1	5.15	122.88	120.30
1	A	294	GLU	CG-CD-OE2	5.11	128.51	118.30
1	A	11	PHE	CB-CG-CD2	-5.10	117.23	120.80
1	A	26	PHE	CB-CG-CD1	-5.09	117.24	120.80
1	A	152	ARG	NE-CZ-NH2	-5.04	117.78	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	178	PHE	O-C-N	5.02	130.73	122.70
1	A	175	ASP	CB-CG-OD2	-5.00	113.80	118.30

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	117	ARG	Sidechain
1	A	177	ARG	Sidechain
1	A	208	ARG	Sidechain
1	A	23	ARG	Sidechain
1	A	321	ARG	Sidechain
1	A	331	ARG	Sidechain
1	A	41	ARG	Sidechain

## 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	3049	0	2914	19	0
2	A	20	0	19	1	0
3	A	2	0	0	0	0
4	A	362	0	0	6	0
All	All	3433	0	2933	19	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 3.

All (19) 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:96:HIS:HD2	1:A:98:VAL:H	1.30	0.80
1:A:92:ASN:HD21	1:A:95:THR:H	1.39	0.70
1:A:93:LEU:HD21	4:A:618:HOH:O	1.93	0.68
1:A:219:GLY:H	1:A:247:ASN:HD21	1.43	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:GLY:HA2	4:A:608:HOH:O	2.01	0.60
1:A:218:VAL:HG11	4:A:608:HOH:O	2.02	0.59
1:A:24:ASP:HB2	1:A:25:PRO:HD2	1.89	0.53
1:A:92:ASN:ND2	1:A:95:THR:H	2.07	0.51
1:A:266:ARG:HG3	4:A:454:HOH:O	2.11	0.51
1:A:119:THR:N	4:A:618:HOH:O	2.43	0.50
1:A:96:HIS:CD2	1:A:98:VAL:H	2.20	0.48
1:A:92:ASN:ND2	1:A:94:PHE:H	2.15	0.45
1:A:337:GLU:HG2	1:A:340:ARG:NH2	2.32	0.44
1:A:91:THR:HG21	4:A:618:HOH:O	2.18	0.43
1:A:137:TRP:HB3	2:A:393[A]:XYS:H51	2.02	0.42
1:A:92:ASN:C	1:A:92:ASN:HD22	2.22	0.42
1:A:323:ASP:HA	1:A:324:PRO:HD3	1.95	0.41
1:A:92:ASN:HD22	1:A:94:PHE:H	1.66	0.41
1:A:381:ASP:HB3	1:A:387[B]:ARG:HD2	2.02	0.41

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	390/387 (101%)	376 (96%)	13 (3%)	1 (0%)	46 23

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	186	GLU

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	306/303 (101%)	297 (97%)	9 (3%)	50 21

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	ARG
1	A	91	THR
1	A	92	ASN
1	A	175	ASP
1	A	185	ASN
1	A	230	HIS
1	A	247	ASN
1	A	347	LEU
1	A	348	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	92	ASN
1	A	96	HIS
1	A	185	ASN
1	A	215	ASN
1	A	222	GLN
1	A	230	HIS
1	A	247	ASN

### 5.3.3 RNA ⓘ

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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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$
2	XLS	A	393[B]	-	9,9,9	1.61	1 (11%)	11,11,11	1.86	2 (18%)
2	XYS	A	393[A]	-	10,10,10	2.03	2 (20%)	12,14,14	2.44	7 (58%)

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	XLS	A	393[B]	-	-	0/10/12/12	0/0/0/0
2	XYS	A	393[A]	-	-	0/0/17/17	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	393[A]	XYS	O5-C5	-4.57	1.36	1.43
2	A	393[A]	XYS	O5-C1	-3.45	1.38	1.43
2	A	393[B]	XLS	O1-C1	4.46	1.40	1.19

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	393[B]	XLS	O1-C1-C2	-4.02	113.89	125.60
2	A	393[A]	XYS	C1-C2-C3	-3.85	104.70	110.43
2	A	393[A]	XYS	O3-C3-C4	2.06	113.72	110.00
2	A	393[A]	XYS	O1-C1-O5	2.24	116.27	109.90
2	A	393[A]	XYS	O4-C4-C5	2.37	113.97	109.21
2	A	393[A]	XYS	O4-C4-C3	2.91	115.98	110.12
2	A	393[A]	XYS	O1-C1-C2	3.11	117.55	109.21
2	A	393[B]	XLS	O2-C2-C1	3.27	118.05	110.22
2	A	393[A]	XYS	O5-C5-C4	4.44	118.06	110.86

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	A	393[A]	XYS	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.