



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2018 – 04:05 AM EST

PDB ID : 4MOH  
Title : Pyranose 2-oxidase V546C mutant with 2-fluorinated glucose  
Authors : Tan, T.C.; Spadiut, O.; Gandini, R.; Haltrich, D.; Divne, C.  
Deposited on : 2013-09-12  
Resolution : 2.10 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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030736  
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-20030736

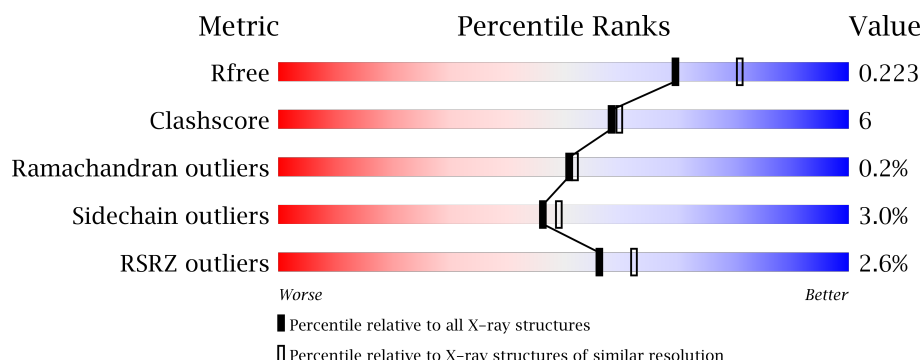
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	4243 (2.10-2.10)
Clashscore	112137	4788 (2.10-2.10)
Ramachandran outliers	110173	4740 (2.10-2.10)
Sidechain outliers	110143	4741 (2.10-2.10)
RSRZ outliers	101464	4275 (2.10-2.10)

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.

Mol	Chain	Length	Quality of chain
1	A	633	<div> <div>2%</div> <div>79%</div> <div>10%</div> <div>9%</div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4956 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.

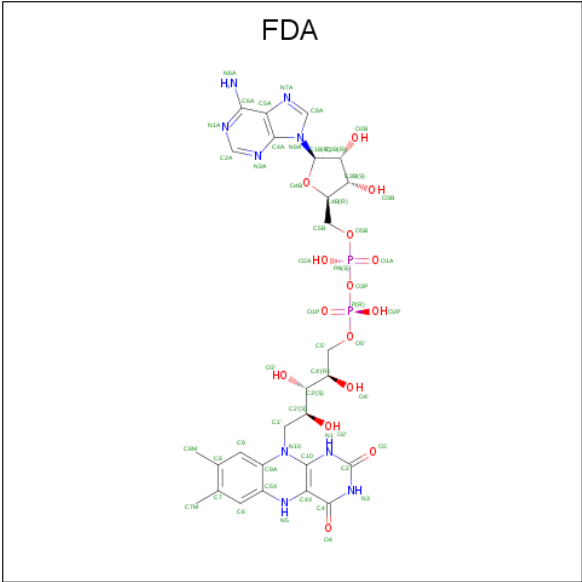
- Molecule 1 is a protein called Pyranose 2-oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	576	Total	C	N	O	S	0	0	0
			4541	2866	777	872	26			

There are 13 discrepancies between the modelled and reference sequences:

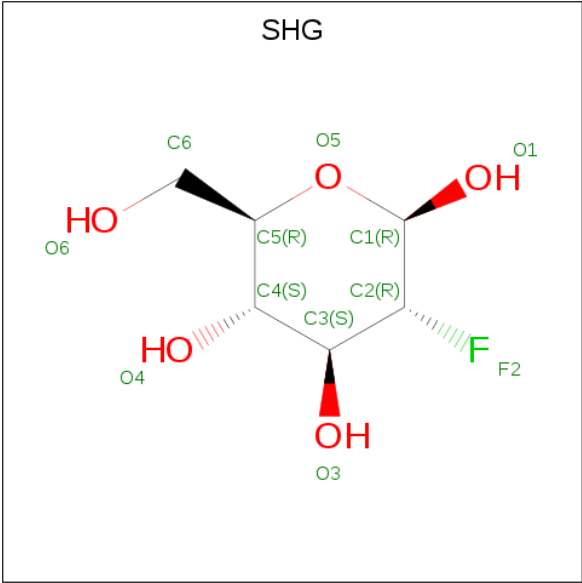
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	ALA	SER	SEE REMARK 999	UNP Q7ZA32
A	546	CYS	VAL	ENGINEERED MUTATION	UNP Q7ZA32
A	623	ALA	-	EXPRESSION TAG	UNP Q7ZA32
A	624	ALA	-	EXPRESSION TAG	UNP Q7ZA32
A	625	ALA	-	EXPRESSION TAG	UNP Q7ZA32
A	626	LEU	-	EXPRESSION TAG	UNP Q7ZA32
A	627	GLU	-	EXPRESSION TAG	UNP Q7ZA32
A	628	HIS	-	EXPRESSION TAG	UNP Q7ZA32
A	629	HIS	-	EXPRESSION TAG	UNP Q7ZA32
A	630	HIS	-	EXPRESSION TAG	UNP Q7ZA32
A	631	HIS	-	EXPRESSION TAG	UNP Q7ZA32
A	632	HIS	-	EXPRESSION TAG	UNP Q7ZA32
A	633	HIS	-	EXPRESSION TAG	UNP Q7ZA32

- Molecule 2 is DIHYDROFLAVINE-ADENINE DINUCLEOTIDE (three-letter code: FDA) (formula: C<sub>27</sub>H<sub>35</sub>N<sub>9</sub>O<sub>15</sub>P<sub>2</sub>).



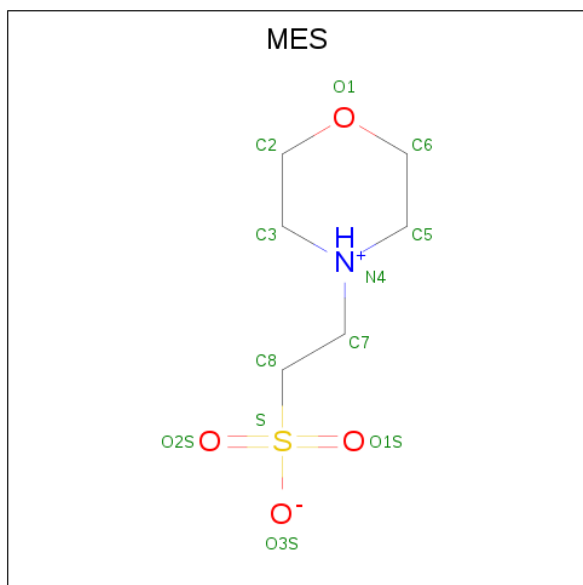
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is 2-deoxy-2-fluoro-beta-D-glucopyranose (three-letter code: SHG) (formula: C<sub>6</sub>H<sub>11</sub>FO<sub>5</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	F	O	0	0
			12	6	1	5		

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C<sub>6</sub>H<sub>13</sub>NO<sub>4</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
4	A	1	12	6	1	4	1	0	0

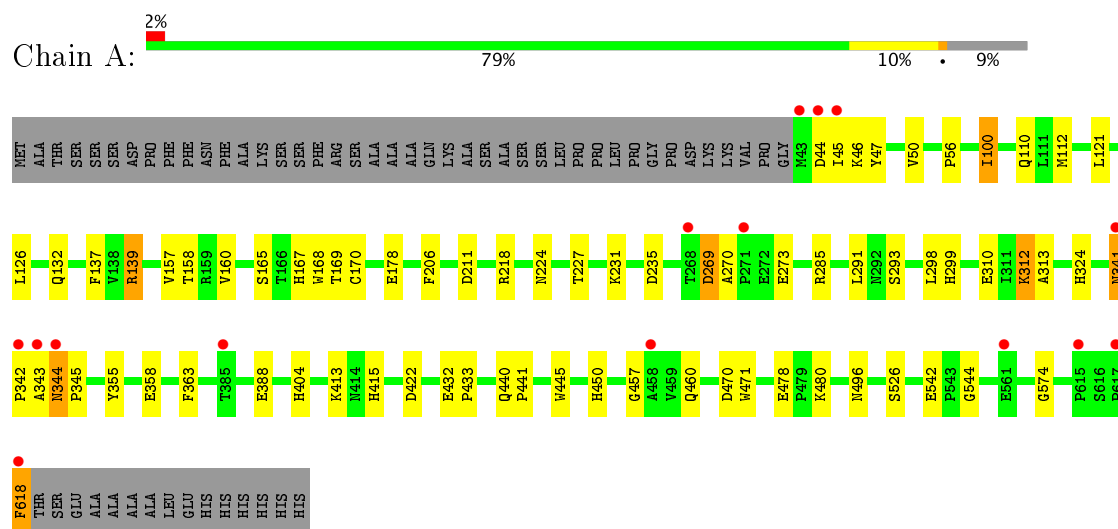
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	338	Total	O	0	0
			338	338		

### 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: Pyranose 2-oxidase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	102.49 Å   102.49 Å   118.10 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	40.00 – 2.10 36.75 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.00-2.10) 99.9 (36.75-2.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.98 (at 2.10 Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, $R_{free}$	0.164   ,   0.221 0.171   ,   0.223	Depositor DCC
$R_{free}$ test set	1163 reflections (3.22%)	DCC
Wilson B-factor (Å <sup>2</sup> )	24.1	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 50.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4956	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FDA, MES, SHG

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	0.95	2/4657 (0.0%)	0.92	6/6331 (0.1%)

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	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	445	TRP	CD2-CE2	5.88	1.48	1.41
1	A	478	GLU	CD-OE1	5.43	1.31	1.25

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	139	ARG	N-CA-CB	-7.07	97.88	110.60
1	A	211	ASP	CB-CG-OD1	6.84	124.46	118.30
1	A	235	ASP	CB-CG-OD2	-6.52	112.43	118.30
1	A	470	ASP	CB-CG-OD1	5.54	123.29	118.30
1	A	139	ARG	CA-CB-CG	5.43	125.35	113.40
1	A	235	ASP	CB-CG-OD1	5.43	123.19	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	A	341	ASN	Peptide

## 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	4541	0	4384	50	0
2	A	53	0	31	0	0
3	A	12	0	10	0	0
4	A	12	0	13	0	0
5	A	338	0	0	4	0
All	All	4956	0	4438	50	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 6.

All (50) 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:344:ASN:H	1:A:345:PRO:HD2	1.04	1.14
1:A:344:ASN:H	1:A:345:PRO:CD	1.61	1.12
1:A:110:GLN:HE21	1:A:167:HIS:HD1	1.10	0.99
1:A:344:ASN:N	1:A:345:PRO:CD	2.29	0.93
1:A:344:ASN:N	1:A:345:PRO:HD2	1.86	0.86
1:A:342:PRO:HD2	1:A:343:ALA:H	1.53	0.73
1:A:343:ALA:O	1:A:344:ASN:HB2	1.90	0.71
1:A:224:ASN:O	1:A:227:THR:HG22	1.94	0.67
1:A:457:GLY:H	1:A:460:GLN:HE21	1.40	0.67
1:A:299:HIS:NE2	1:A:310:GLU:OE2	2.31	0.64
1:A:45:ILE:HG22	1:A:312:LYS:HE2	1.83	0.60
1:A:46:LYS:HD3	1:A:312:LYS:HG3	1.84	0.60
1:A:343:ALA:O	1:A:344:ASN:CB	2.50	0.59
1:A:404:HIS:HE1	5:A:1083:HOH:O	1.87	0.58
1:A:100:ILE:HD13	1:A:100:ILE:C	2.24	0.58
1:A:285:ARG:O	1:A:298:LEU:HD12	2.03	0.57
1:A:178:GLU:OE2	1:A:441:PRO:HD3	2.05	0.57
1:A:342:PRO:CD	1:A:343:ALA:H	2.18	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:ASN:HB3	1:A:342:PRO:HD3	1.87	0.55
1:A:50:VAL:HG13	1:A:313:ALA:HB2	1.87	0.55
1:A:218:ARG:HD2	5:A:908:HOH:O	2.08	0.54
1:A:270:ALA:HB1	1:A:273:GLU:HG3	1.88	0.54
1:A:45:ILE:HG22	1:A:312:LYS:CE	2.38	0.54
1:A:415:HIS:NE2	1:A:422:ASP:OD2	2.35	0.53
1:A:45:ILE:CG2	1:A:312:LYS:HE2	2.43	0.49
1:A:158:THR:HG22	1:A:160:VAL:HG22	1.95	0.48
1:A:355:TYR:HA	1:A:480:LYS:O	2.14	0.48
1:A:344:ASN:N	1:A:345:PRO:HD3	2.25	0.48
1:A:47:TYR:O	1:A:313:ALA:HA	2.14	0.47
1:A:618:PHE:CD1	1:A:618:PHE:C	2.87	0.47
1:A:342:PRO:CD	1:A:343:ALA:N	2.78	0.47
1:A:358:GLU:HG2	1:A:544:GLY:HA2	1.97	0.46
1:A:342:PRO:HD2	1:A:343:ALA:N	2.28	0.46
1:A:137:PHE:O	1:A:139:ARG:HB2	2.16	0.46
1:A:471:TRP:CH2	1:A:526:SER:HA	2.52	0.45
1:A:224:ASN:HA	1:A:227:THR:HG22	1.98	0.45
1:A:293:SER:HA	1:A:574:GLY:O	2.18	0.44
1:A:440:GLN:NE2	5:A:1187:HOH:O	2.51	0.44
1:A:291:LEU:HD23	1:A:291:LEU:HA	1.88	0.44
1:A:269:ASP:N	1:A:269:ASP:OD2	2.51	0.43
1:A:169:THR:O	1:A:170:CYS:HB2	2.19	0.43
1:A:618:PHE:C	1:A:618:PHE:HD1	2.21	0.42
1:A:270:ALA:HB1	1:A:273:GLU:CG	2.49	0.42
1:A:56:PRO:HD3	1:A:165:SER:HB3	2.01	0.42
1:A:45:ILE:HB	5:A:1098:HOH:O	2.19	0.42
1:A:363:PHE:HA	1:A:471:TRP:O	2.20	0.41
1:A:432:GLU:HB2	1:A:433:PRO:HD2	2.03	0.40
1:A:167:HIS:CD2	1:A:167:HIS:C	2.94	0.40
1:A:126:LEU:HD12	1:A:132:GLN:HG3	2.02	0.40
1:A:157:VAL:HG21	1:A:324:HIS:HE1	1.87	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	574/633 (91%)	556 (97%)	17 (3%)	1 (0%)	51 52

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	344	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	504/548 (92%)	489 (97%)	15 (3%)	46 49

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

Mol	Chain	Res	Type
1	A	44	ASP
1	A	100	ILE
1	A	112	MET
1	A	121	LEU
1	A	168	TRP
1	A	206	PHE
1	A	231	LYS
1	A	269	ASP
1	A	312	LYS
1	A	388	GLU
1	A	413	LYS
1	A	450	HIS
1	A	496	ASN
1	A	542	GLU
1	A	618	PHE

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

Mol	Chain	Res	Type
1	A	404	HIS
1	A	460	GLN

### 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 ⓘ

3 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	FDA	A	801	1	51,58,58	1.47	7 (13%)	54,89,89	3.31	14 (25%)
3	SHG	A	802	-	11,12,12	0.83	0	14,17,17	3.91	8 (57%)
4	MES	A	803	-	12,12,12	1.99	2 (16%)	14,16,16	2.97	4 (28%)

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	FDA	A	801	1	-	0/28/50/50	0/6/6/6
3	SHG	A	802	-	-	0/2/22/22	0/1/1/1
4	MES	A	803	-	-	0/6/14/14	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	803	MES	C8-S	-5.81	1.68	1.77
2	A	801	FDA	O3B-C3B	-2.13	1.38	1.43
2	A	801	FDA	C4-N3	2.02	1.36	1.33
4	A	803	MES	C3-N4	2.05	1.52	1.47
2	A	801	FDA	C8M-C8	2.53	1.56	1.51
2	A	801	FDA	C9A-N10	2.79	1.42	1.38
2	A	801	FDA	C2A-N3A	2.92	1.37	1.32
2	A	801	FDA	C1'-N10	3.91	1.52	1.48
2	A	801	FDA	C4-C4X	5.50	1.51	1.41

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	FDA	C4X-C4-N3	-10.47	108.58	123.48
2	A	801	FDA	N3A-C2A-N1A	-10.06	120.10	128.86
4	A	803	MES	O1S-S-C8	-2.88	104.31	106.79
4	A	803	MES	O1-C2-C3	-2.51	106.22	111.83
2	A	801	FDA	C4A-C5A-N7A	-2.11	107.37	109.41
3	A	802	SHG	O4-C4-C3	2.04	114.79	110.36
4	A	803	MES	C6-C5-N4	2.31	113.34	110.11
3	A	802	SHG	C3-C4-C5	2.41	114.47	110.22
3	A	802	SHG	O5-C1-C2	2.70	119.42	110.52
2	A	801	FDA	C2B-C3B-C4B	2.70	107.88	102.62
2	A	801	FDA	O2P-P-O1P	2.73	126.42	112.28
2	A	801	FDA	C5A-C6A-N6A	2.95	126.47	120.47
2	A	801	FDA	O4B-C4B-C5B	3.08	119.81	109.40
2	A	801	FDA	C2A-N1A-C6A	3.14	124.26	118.77
2	A	801	FDA	C5B-C4B-C3B	3.17	127.38	115.29
2	A	801	FDA	O2B-C2B-C3B	3.50	123.05	111.83
3	A	802	SHG	C6-C5-C4	3.65	121.55	113.00
3	A	802	SHG	O5-C5-C4	3.68	116.45	109.66
2	A	801	FDA	O2B-C2B-C1B	3.81	123.54	111.61
3	A	802	SHG	O3-C3-C4	4.07	119.20	110.36
2	A	801	FDA	C4X-N5-C5X	4.09	121.08	116.76
2	A	801	FDA	C1'-N10-C9A	5.29	123.19	118.35
3	A	802	SHG	C1-O5-C5	6.30	124.75	113.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	803	MES	O2S-S-C8	9.59	115.03	106.79
3	A	802	SHG	F2-C2-C3	10.56	116.23	108.49
2	A	801	FDA	C4-N3-C2	15.11	128.37	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	576/633 (90%)	-0.30	15 (2%) 56 62	15, 25, 44, 92	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	342	PRO	7.3
1	A	341	ASN	6.8
1	A	343	ALA	5.7
1	A	45	ILE	5.3
1	A	44	ASP	4.9
1	A	618	PHE	4.6
1	A	344	ASN	4.5
1	A	43	MET	3.7
1	A	617	PRO	2.9
1	A	268	THR	2.7
1	A	615	PRO	2.6
1	A	385	THR	2.5
1	A	561	GLU	2.3
1	A	271	PRO	2.2
1	A	458	ALA	2.2

### 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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	SHG	A	802	12/12	0.91	0.12	0.57	33,37,40,40	0
4	MES	A	803	12/12	0.97	0.08	0.32	22,24,25,26	0
2	FDA	A	801	53/53	0.97	0.11	-0.35	17,20,23,23	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.