



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 04:52 am GMT

PDB ID : 4WVC  
Title : Crystal structure of GH63 mannosylglycerate hydrolase from *Thermus thermophilus* HB8 in complex with Tris and D-glycerate  
Authors : Miyazaki, T.; Ichikawa, M.; Nishikawa, A.; Tono-zuka, T.  
Deposited on : 2014-11-05  
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 : trunk28620  
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) : recalc28949

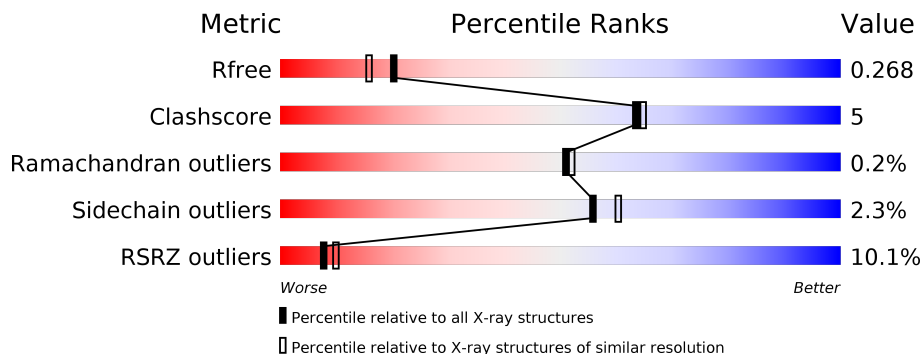
# 1 Overall quality at a glance

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	420	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="width: 89%; height: 10px; background-color: green;"></div> <div style="width: 10%; height: 10px; background-color: yellow;"></div> <div style="width: 1%; height: 10px; background-color: grey;"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> <span>%</span> <span>89%</span> <span>10%</span> <span>.</span> </div> </div>
1	B	420	<div> <div style="width: 19%; height: 10px; background-color: red;"></div> <div style="width: 83%; height: 10px; background-color: green;"></div> <div style="width: 13%; height: 10px; background-color: yellow;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> <span>19%</span> <span>83%</span> <span>13%</span> <span>..</span> </div> </div>

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	TRS	A	501	-	X	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7093 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.

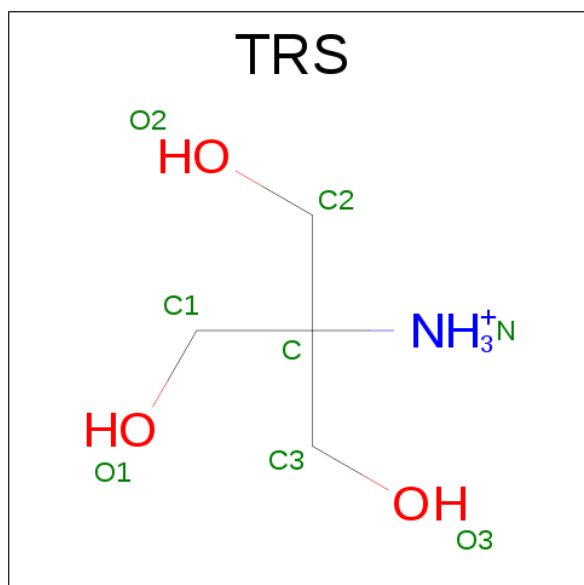
- Molecule 1 is a protein called Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	413	Total	C	N	O	S	0	0	0
			3416	2215	598	598	5			
1	B	409	Total	C	N	O	S	0	0	0
			3380	2191	590	594	5			

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

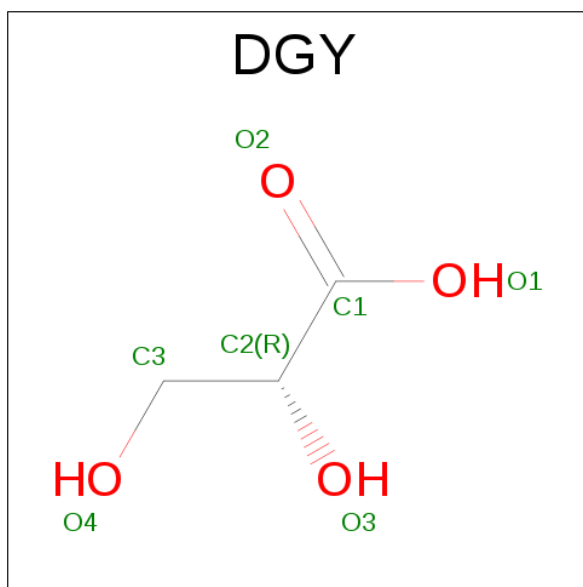
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		
2	A	1	Total	Cl	0	0
			1	1		

- Molecule 3 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: C<sub>4</sub>H<sub>12</sub>NO<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			8	4	1	3		
3	B	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 4 is (2R)-2,3-DIHYDROXYPROPANOIC ACID (three-letter code: DGY) (formula: C<sub>3</sub>H<sub>6</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	3	4		
4	B	1	Total	C	O	0	0
			7	3	4		

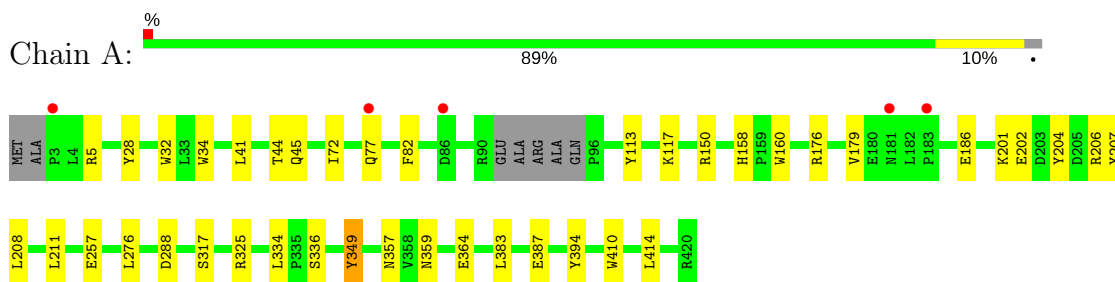
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	145	Total	O	0	0
			145	145		
5	B	120	Total	O	0	0
			120	120		

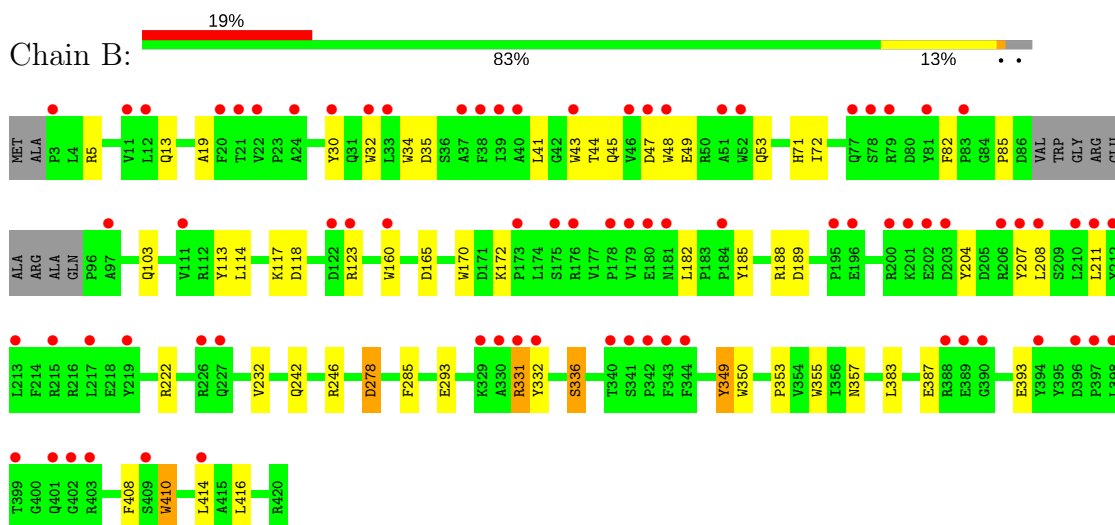
### 3 Residue-property plots [i](#)

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: Uncharacterized protein



- Molecule 1: Uncharacterized protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	92.94Å 92.94Å 190.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.15 – 2.10 45.14 – 1.96	Depositor EDS
% Data completeness (in resolution range)	97.5 (45.15-2.10) 98.2 (45.14-1.96)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.67 (at 1.97Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, $R_{free}$	0.245 , 0.280 0.256 , 0.268	Depositor DCC
$R_{free}$ test set	2859 reflections (5.39%)	DCC
Wilson B-factor (Å <sup>2</sup> )	17.2	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 52.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.054 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	7093	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.57% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: TRS, DGY, CL

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	0.54	0/3535	0.71	2/4818 (0.0%)
1	B	0.53	0/3497	0.71	2/4766 (0.0%)
All	All	0.53	0/7032	0.71	4/9584 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	278	ASP	CB-CG-OD1	6.88	124.50	118.30
1	A	325	ARG	NE-CZ-NH2	-5.58	117.51	120.30
1	A	325	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	B	278	ASP	CB-CG-OD2	-5.29	113.54	118.30

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	3416	0	3280	23	0
1	B	3380	0	3245	45	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	8	0	12	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	8	0	12	0	0
4	A	7	0	5	0	0
4	B	7	0	5	2	0
5	A	145	0	0	4	0
5	B	120	0	0	14	1
All	All	7093	0	6559	66	1

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 (66) 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:257:GLU:OE1	1:B:222:ARG:NE	1.97	0.95
1:A:201:LYS:HG3	5:A:711:HOH:O	1.89	0.73
1:A:364:GLU:HB3	5:A:635:HOH:O	1.90	0.72
1:B:47:ASP:HB3	5:B:652:HOH:O	1.89	0.71
1:B:165:ASP:HB3	5:B:705:HOH:O	1.92	0.68
1:B:43:TRP:HE3	5:B:652:HOH:O	1.77	0.68
1:B:19:ALA:HA	5:B:623:HOH:O	1.95	0.66
1:B:53:GLN:HB3	5:B:623:HOH:O	1.96	0.66
1:A:334:LEU:H	1:A:359:ASN:HD21	1.44	0.64
1:B:336:SER:OG	1:B:353:PRO:O	2.14	0.64
1:B:5:ARG:HB3	5:B:672:HOH:O	1.98	0.62
1:B:82:PHE:HE2	1:B:208:LEU:HD11	1.65	0.61
1:A:158:HIS:HD2	1:A:160:TRP:H	1.49	0.61
1:B:242:GLN:HE21	1:B:246:ARG:HE	1.49	0.60
1:A:257:GLU:OE1	1:B:222:ARG:CZ	2.53	0.57
1:B:30:TYR:HD1	5:B:660:HOH:O	1.87	0.56
1:A:82:PHE:HE1	1:A:208:LEU:HD11	1.69	0.56
1:B:211:LEU:O	1:B:211:LEU:HD23	2.06	0.56
1:B:182:LEU:HB3	5:B:704:HOH:O	2.07	0.54
1:B:48:TRP:CH2	1:B:123:ARG:HG3	2.43	0.54
1:A:364:GLU:HG2	5:A:732:HOH:O	2.08	0.52
1:B:13:GLN:HB3	5:B:620:HOH:O	2.09	0.52
1:B:189:ASP:HA	5:B:707:HOH:O	2.09	0.51
1:B:82:PHE:HB2	1:B:188:ARG:HD2	1.93	0.51
1:B:383:LEU:O	1:B:387:GLU:HG2	2.11	0.51
1:B:331:ARG:HG3	1:B:332:TYR:CD2	2.46	0.51
1:B:357:ASN:HA	1:B:414:LEU:CD1	2.41	0.51
1:B:5:ARG:C	5:B:672:HOH:O	2.50	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:179:VAL:HG11	1:A:206:ARG:HD2	1.93	0.50
1:A:211:LEU:HD23	1:A:211:LEU:O	2.12	0.50
1:B:331:ARG:C	1:B:331:ARG:CD	2.80	0.48
1:A:202:GLU:O	1:A:206:ARG:HD3	2.13	0.48
1:A:5:ARG:HH22	1:A:45:GLN:HE21	1.62	0.48
1:A:82:PHE:CE1	1:A:208:LEU:HD11	2.48	0.48
1:B:278:ASP:OD2	1:B:285:PHE:HE2	1.97	0.48
1:B:5:ARG:CB	5:B:672:HOH:O	2.59	0.48
1:B:211:LEU:HD23	1:B:211:LEU:C	2.34	0.47
1:B:82:PHE:CZ	1:B:185:TYR:HB2	2.49	0.47
1:A:176:ARG:NH2	1:A:288:ASP:OD2	2.47	0.47
1:B:41:LEU:O	1:B:44:THR:OG1	2.20	0.47
1:A:383:LEU:O	1:A:387:GLU:HG2	2.15	0.46
1:A:211:LEU:HD23	1:A:211:LEU:C	2.36	0.45
1:B:160:TRP:CD1	4:B:502:DGY:H3C1	2.51	0.45
1:A:204:TYR:HA	1:A:207:TYR:CD2	2.51	0.45
1:B:350:TRP:CE2	1:B:355:TRP:CZ2	3.05	0.45
1:B:113:TYR:CE1	1:B:117:LYS:HG3	2.51	0.45
1:B:117:LYS:O	1:B:118:ASP:C	2.56	0.44
1:B:5:ARG:HH22	1:B:45:GLN:HE21	1.66	0.44
1:A:357:ASN:HA	1:A:414:LEU:CD1	2.48	0.43
1:B:165:ASP:CB	5:B:705:HOH:O	2.57	0.43
1:B:204:TYR:HA	1:B:207:TYR:CD2	2.52	0.43
1:A:113:TYR:CE1	1:A:117:LYS:HG3	2.54	0.43
1:A:41:LEU:O	1:A:44:THR:OG1	2.21	0.42
1:B:47:ASP:HA	5:B:674:HOH:O	2.18	0.42
1:B:393:GLU:HG3	1:B:408:PHE:CE1	2.54	0.42
1:B:172:LYS:NZ	1:B:293:GLU:OE1	2.53	0.42
1:B:170:TRP:CZ3	1:B:232:VAL:HG11	2.55	0.42
1:B:45:GLN:HB2	1:B:416:LEU:HD23	2.01	0.42
1:B:48:TRP:CE3	1:B:49:GLU:N	2.89	0.41
1:A:77:GLN:NE2	5:A:724:HOH:O	2.54	0.41
1:B:71:HIS:HB2	1:B:103:GLN:HG3	2.01	0.41
1:B:35:ASP:HA	1:B:410:TRP:HE3	1.84	0.41
1:B:32:TRP:CE2	1:B:72:ILE:HG12	2.56	0.40
1:A:32:TRP:CE2	1:A:72:ILE:HG12	2.57	0.40
1:B:165:ASP:OD2	4:B:502:DGY:O3	2.30	0.40
1:A:28:TYR:CZ	3:A:501:TRS:H21	2.56	0.40

All (1) 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)
5:B:633:HOH:O	5:B:633:HOH:O[2_655]	0.70	1.50

## 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	409/420 (97%)	395 (97%)	13 (3%)	1 (0%)	51	52
1	B	405/420 (96%)	390 (96%)	14 (4%)	1 (0%)	51	52
All	All	814/840 (97%)	785 (96%)	27 (3%)	2 (0%)	51	52

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	349	TYR
1	A	349	TYR

### 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	345/349 (99%)	336 (97%)	9 (3%)	51	55
1	B	342/349 (98%)	335 (98%)	7 (2%)	60	66
All	All	687/698 (98%)	671 (98%)	16 (2%)	56	60

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	TRP
1	A	150	ARG
1	A	186	GLU
1	A	276	LEU
1	A	317	SER
1	A	336	SER
1	A	349	TYR
1	A	394	TYR
1	A	410	TRP
1	B	34	TRP
1	B	85	PRO
1	B	114	LEU
1	B	331	ARG
1	B	336	SER
1	B	349	TYR
1	B	410	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	GLN
1	A	75	HIS
1	A	77	GLN
1	A	158	HIS
1	A	227	GLN
1	A	359	ASN
1	B	15	ASN
1	B	45	GLN
1	B	75	HIS
1	B	192	HIS
1	B	242	GLN
1	B	245	ASN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	TRS	A	501	-	7,7,7	0.98	0	9,9,9	4.48	7 (77%)
4	DGY	A	502	-	2,6,6	0.47	0	1,7,7	0.80	0
3	TRS	B	501	-	7,7,7	0.53	0	9,9,9	0.22	0
4	DGY	B	502	-	2,6,6	0.50	0	1,7,7	0.04	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	TRS	A	501	-	-	0/9/9/9	0/0/0/0
4	DGY	A	502	-	-	0/2/6/6	0/0/0/0
3	TRS	B	501	-	-	0/9/9/9	0/0/0/0
4	DGY	B	502	-	-	0/2/6/6	0/0/0/0

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	TRS	C2-C-N	-7.06	92.72	107.73
3	A	501	TRS	C1-C-N	-6.09	94.77	107.73
3	A	501	TRS	C3-C-N	-5.99	94.99	107.73
3	A	501	TRS	O1-C1-C	-3.68	99.82	110.47
3	A	501	TRS	O2-C2-C	-2.75	102.50	110.47
3	A	501	TRS	C2-C-C1	4.09	122.65	111.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	TRS	C3-C-C2	4.12	122.74	111.06

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	TRS	1	0
4	B	502	DGY	2	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	413/420 (98%)	0.08	5 (1%) 79 82	14, 23, 39, 54	0
1	B	409/420 (97%)	1.09	78 (19%) 1 1	18, 35, 54, 69	0
All	All	822/840 (97%)	0.58	83 (10%) 8 10	14, 28, 49, 69	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	48	TRP	6.7
1	B	212	TYR	5.8
1	B	398	LEU	4.7
1	B	180	GLU	4.4
1	B	181	ASN	4.0
1	B	219	TYR	4.0
1	B	401	GLN	3.8
1	B	402	GLY	3.8
1	B	12	LEU	3.8
1	B	160	TRP	3.7
1	B	77	GLN	3.7
1	A	77	GLN	3.6
1	B	329	LYS	3.6
1	B	184	PRO	3.4
1	B	226	ARG	3.4
1	B	3	PRO	3.2
1	B	331	ARG	3.2
1	B	51	ALA	2.9
1	B	11	VAL	2.9
1	B	388	ARG	2.9
1	B	176	ARG	2.9
1	B	123	ARG	2.8
1	B	341	SER	2.8
1	B	78	SER	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	178	PRO	2.8
1	B	397	PRO	2.8
1	B	30	TYR	2.8
1	A	181	ASN	2.7
1	B	37	ALA	2.7
1	B	227	GLN	2.7
1	B	215	ARG	2.7
1	B	394	TYR	2.7
1	B	39	ILE	2.6
1	B	414	LEU	2.6
1	B	175	SER	2.6
1	B	79	ARG	2.6
1	B	97	ALA	2.6
1	B	203	ASP	2.6
1	B	217	LEU	2.6
1	B	340	THR	2.6
1	B	403	ARG	2.5
1	B	43	TRP	2.5
1	B	342	PRO	2.4
1	A	3	PRO	2.4
1	A	86	ASP	2.4
1	B	332	TYR	2.4
1	B	399	THR	2.4
1	B	22	VAL	2.4
1	B	330	ALA	2.4
1	B	32	TRP	2.4
1	B	344	PHE	2.4
1	B	179	VAL	2.4
1	A	183	PRO	2.3
1	B	195	PRO	2.3
1	B	46	VAL	2.3
1	B	47	ASP	2.3
1	B	81	TYR	2.2
1	B	210	LEU	2.2
1	B	20	PHE	2.2
1	B	38	PHE	2.2
1	B	206	ARG	2.2
1	B	122	ASP	2.2
1	B	202	GLU	2.2
1	B	24	ALA	2.2
1	B	390	GLY	2.2
1	B	83	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	389	GLU	2.2
1	B	111	VAL	2.1
1	B	21	THR	2.1
1	B	208	LEU	2.1
1	B	201	LYS	2.1
1	B	343	PHE	2.1
1	B	196	GLU	2.1
1	B	33	LEU	2.1
1	B	213	LEU	2.1
1	B	211	LEU	2.0
1	B	207	TYR	2.0
1	B	173	PRO	2.0
1	B	396	ASP	2.0
1	B	409	SER	2.0
1	B	52	TRP	2.0
1	B	40	ALA	2.0
1	B	200	ARG	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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(Å <sup>2</sup> )	Q<0.9
3	TRS	A	501	8/8	0.84	0.23	6.20	32,36,38,40	0
3	TRS	B	501	8/8	0.79	0.20	0.98	30,35,38,40	0
4	DGY	B	502	7/7	0.88	0.20	-0.04	28,31,35,36	0
4	DGY	A	502	7/7	0.94	0.11	-0.86	19,19,21,21	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	CL	B	500	1/1	1.00	0.10	-	20,20,20,20	1
2	CL	A	500	1/1	0.99	0.10	-	23,23,23,23	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.