



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 08:43 pm BST

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.; Tonozuka, 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

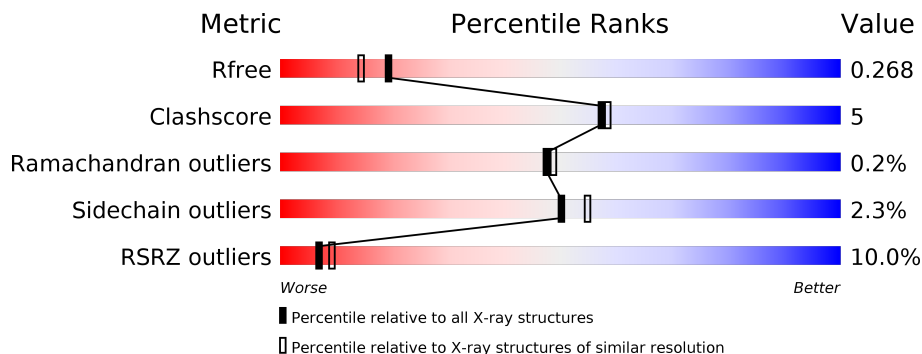
1 Overall quality at a glance

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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (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 ≥ 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 $\leq 5\%$. 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	 18% 89% 10% .
1	B	420	 18% 83% 13% . .

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	-	-

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.

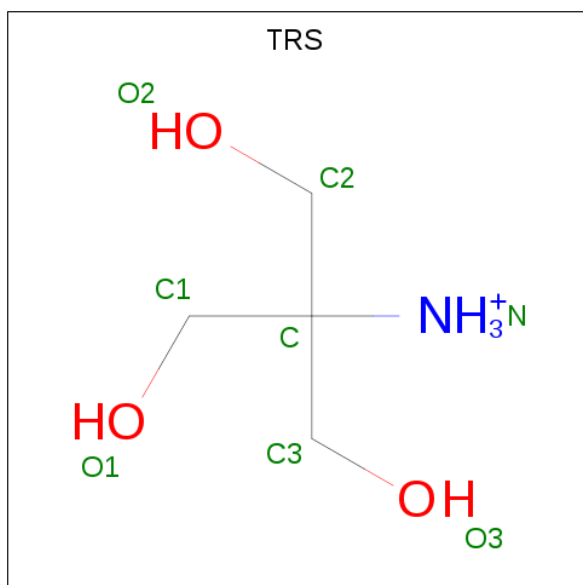
- Molecule 1 is a protein called Uncharacterized protein.

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

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

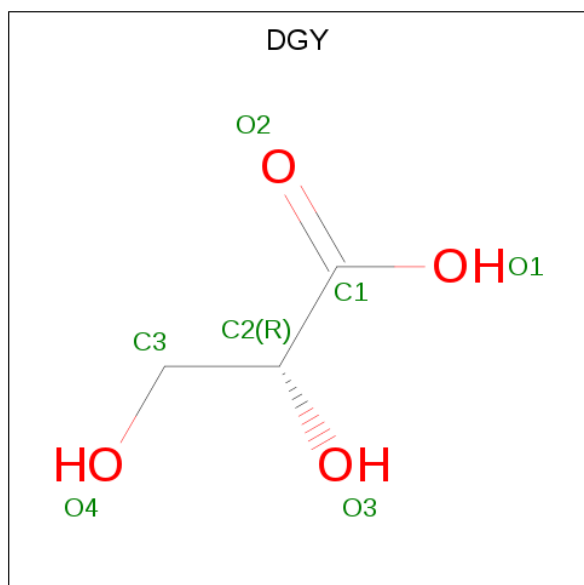
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total 1	Cl 1	0	0
2	A	1	Total 1	Cl 1	0	0

- Molecule 3 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: C₄H₁₂NO₃).



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: C₃H₆O₄).



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		

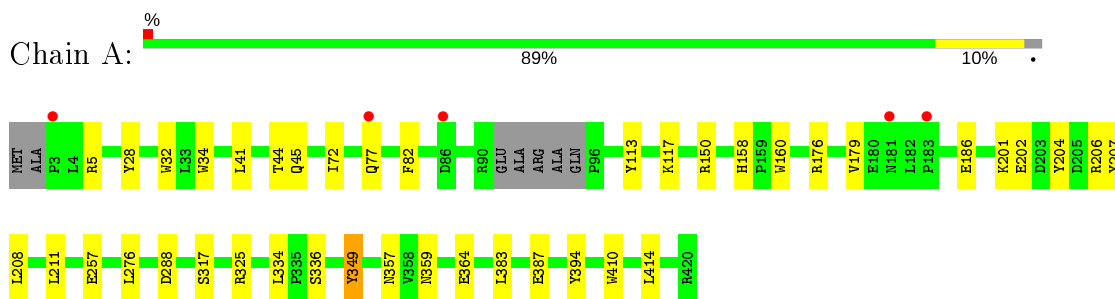
- Molecule 5 is water.

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

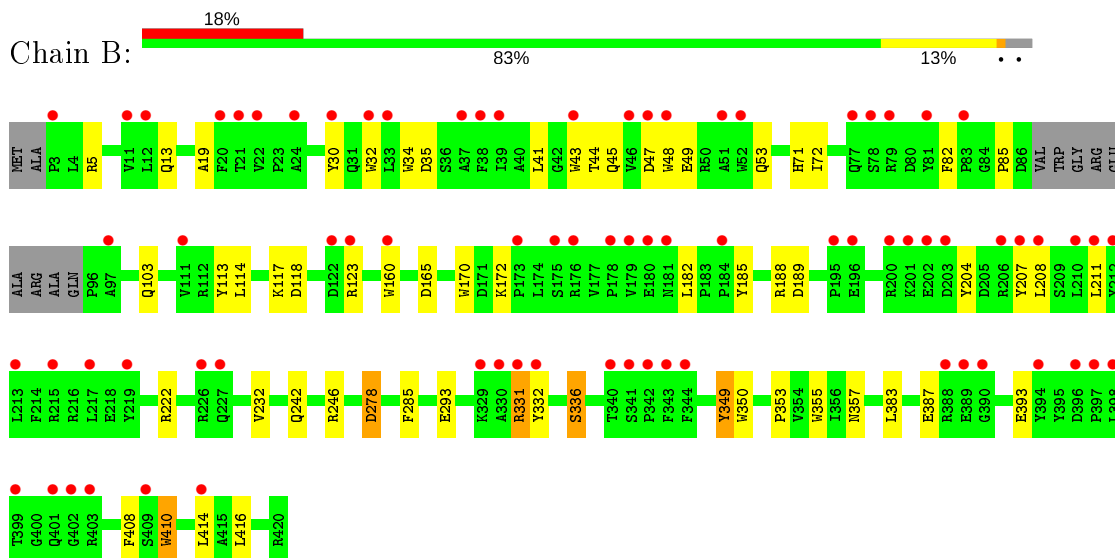
3 Residue-property plots [i](#)

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, α , β , γ	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
$\langle I/\sigma(I) \rangle$ ¹	4.67 (at 1.97Å)	Xtrriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.245 , 0.280 0.256 , 0.268	Depositor DCC
R_{free} test set	3437 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	17.2	Xtrriage
Anisotropy	0.012	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 52.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtrriage
Estimated twinning fraction	0.054 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	7093	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtrriage'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.*

¹Intensities estimated from amplitudes.

²Theoretical 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%)	47	49
1	B	405/420 (96%)	390 (96%)	14 (4%)	1 (0%)	47	49
All	All	814/840 (97%)	785 (96%)	27 (3%)	2 (0%)	47	49

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%)	46	50
1	B	342/349 (98%)	335 (98%)	7 (2%)	55	60
All	All	687/698 (98%)	671 (98%)	16 (2%)	50	55

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	B	501	-	7,7,7	0.46	0	9,9,9	0.17	0
4	DGY	A	502	-	2,6,6	0.47	0	2,7,7	0.84	0
4	DGY	B	502	-	2,6,6	0.51	0	2,7,7	0.11	0
3	TRS	A	501	-	7,7,7	0.80	0	9,9,9	3.61	7 (77%)

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	B	501	-	-	6/9/9/9	-
4	DGY	A	502	-	-	1/2/6/6	-
4	DGY	B	502	-	-	1/2/6/6	-
3	TRS	A	501	-	-	4/9/9/9	-

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	-5.11	92.72	107.98
3	A	501	TRS	C1-C-N	-4.42	94.77	107.98
3	A	501	TRS	C3-C-N	-4.35	94.99	107.98
3	A	501	TRS	C3-C-C2	3.85	122.74	110.81
3	A	501	TRS	C2-C-C1	3.82	122.65	110.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	TRS	O1-C1-C	-3.53	99.82	111.00
3	A	501	TRS	O2-C2-C	-2.68	102.50	111.00

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	TRS	C2-C-C1-O1
3	A	501	TRS	C2-C-C3-O3
3	B	501	TRS	C1-C-C2-O2
3	B	501	TRS	N-C-C2-O2
3	B	501	TRS	N-C-C3-O3
3	A	501	TRS	N-C-C1-O1
3	A	501	TRS	C1-C-C2-O2
4	A	502	DGY	O3-C2-C3-O4
3	B	501	TRS	C3-C-C2-O2
3	B	501	TRS	C1-C-C3-O3
3	B	501	TRS	C2-C-C3-O3
4	B	502	DGY	C1-C2-C3-O4

There are no ring outliers.

2 monomers are involved in 3 short contacts:

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

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	413/420 (98%)	0.08	5 (1%) 79 82	14, 23, 39, 54	0
1	B	409/420 (97%)	1.09	77 (18%) 1 1	18, 35, 54, 69	0
All	All	822/840 (97%)	0.58	82 (9%) 7 9	14, 28, 49, 69	0

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	48	TRP	6.7
1	B	212	TYR	5.9
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	226	ARG	3.4
1	B	184	PRO	3.4
1	B	3	PRO	3.2
1	B	331	ARG	3.2
1	B	11	VAL	2.9
1	B	176	ARG	2.9
1	B	388	ARG	2.9
1	B	51	ALA	2.9
1	B	123	ARG	2.8
1	B	78	SER	2.8
1	B	341	SER	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	178	PRO	2.8
1	B	30	TYR	2.8
1	B	397	PRO	2.8
1	B	227	GLN	2.7
1	A	181	ASN	2.7
1	B	37	ALA	2.7
1	B	215	ARG	2.7
1	B	394	TYR	2.7
1	B	39	ILE	2.6
1	B	175	SER	2.6
1	B	203	ASP	2.6
1	B	217	LEU	2.6
1	B	414	LEU	2.6
1	B	79	ARG	2.6
1	B	97	ALA	2.6
1	B	340	THR	2.6
1	B	403	ARG	2.5
1	B	43	TRP	2.5
1	B	342	PRO	2.5
1	A	3	PRO	2.4
1	B	22	VAL	2.4
1	B	332	TYR	2.4
1	B	399	THR	2.4
1	A	86	ASP	2.4
1	A	183	PRO	2.4
1	B	32	TRP	2.4
1	B	344	PHE	2.4
1	B	179	VAL	2.4
1	B	330	ALA	2.4
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	122	ASP	2.2
1	B	202	GLU	2.2
1	B	206	ARG	2.2
1	B	24	ALA	2.2
1	B	83	PRO	2.2
1	B	390	GLY	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	201	LYS	2.1
1	B	208	LEU	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	52	TRP	2.0
1	B	396	ASP	2.0
1	B	409	SER	2.0
1	B	173	PRO	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. 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	TRS	B	501	8/8	0.79	0.20	30,35,38,40	0
3	TRS	A	501	8/8	0.84	0.23	32,36,38,40	0
4	DGY	B	502	7/7	0.88	0.20	28,31,35,36	0
4	DGY	A	502	7/7	0.94	0.11	19,19,21,21	0
2	CL	A	500	1/1	0.99	0.10	23,23,23,23	1
2	CL	B	500	1/1	1.00	0.10	20,20,20,20	1

6.5 Other polymers

There are no such residues in this entry.