



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2024 – 04:12 PM EDT

PDB ID : 4PA5
Title : Tgl - a bacterial spore coat transglutaminase - cystamine complex
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Deposited on : 2014-04-07
Resolution : 1.86 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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 ⓘ](#)) 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 : 2.36.2
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) : 2.36.2

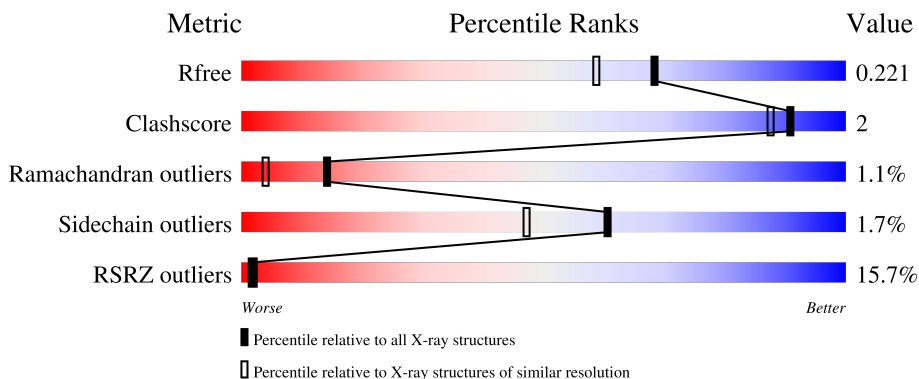
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.86 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	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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 $\leq 5\%$. 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	260	
1	B	260	

2 Entry composition i

There are 10 unique types of molecules in this entry. The entry contains 4269 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 Protein-glutamine gamma-glutamyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	242	1993	1300	333	352	8	0	2	0
1	B	242	1997	1302	331	355	9	0	2	0

There are 30 discrepancies between the modelled and reference sequences:

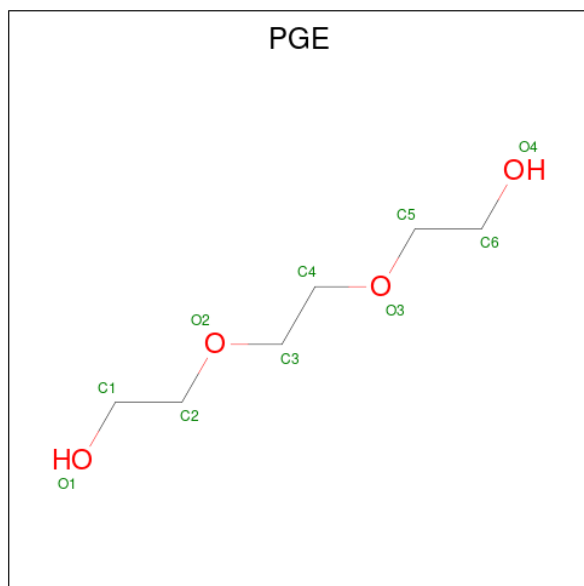
Chain	Residue	Modelled	Actual	Comment	Reference
A	246	VAL	-	expression tag	UNP P40746
A	247	ASP	-	expression tag	UNP P40746
A	248	LYS	-	expression tag	UNP P40746
A	249	LEU	-	expression tag	UNP P40746
A	250	ALA	-	expression tag	UNP P40746
A	251	ALA	-	expression tag	UNP P40746
A	252	ALA	-	expression tag	UNP P40746
A	253	LEU	-	expression tag	UNP P40746
A	254	GLU	-	expression tag	UNP P40746
A	255	HIS	-	expression tag	UNP P40746
A	256	HIS	-	expression tag	UNP P40746
A	257	HIS	-	expression tag	UNP P40746
A	258	HIS	-	expression tag	UNP P40746
A	259	HIS	-	expression tag	UNP P40746
A	260	HIS	-	expression tag	UNP P40746
B	246	VAL	-	expression tag	UNP P40746
B	247	ASP	-	expression tag	UNP P40746
B	248	LYS	-	expression tag	UNP P40746
B	249	LEU	-	expression tag	UNP P40746
B	250	ALA	-	expression tag	UNP P40746
B	251	ALA	-	expression tag	UNP P40746
B	252	ALA	-	expression tag	UNP P40746
B	253	LEU	-	expression tag	UNP P40746
B	254	GLU	-	expression tag	UNP P40746
B	255	HIS	-	expression tag	UNP P40746

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Chain	Residue	Modelled	Actual	Comment	Reference
B	256	HIS	-	expression tag	UNP P40746
B	257	HIS	-	expression tag	UNP P40746
B	258	HIS	-	expression tag	UNP P40746
B	259	HIS	-	expression tag	UNP P40746
B	260	HIS	-	expression tag	UNP P40746

- Molecule 2 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			10	6	4		
2	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

- Molecule 4 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).

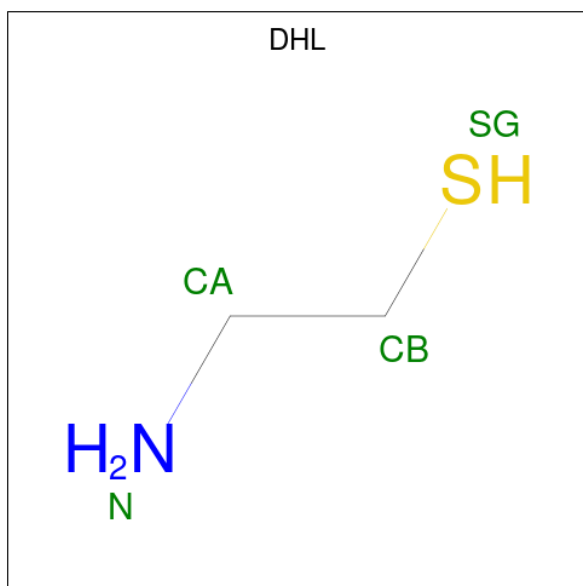


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

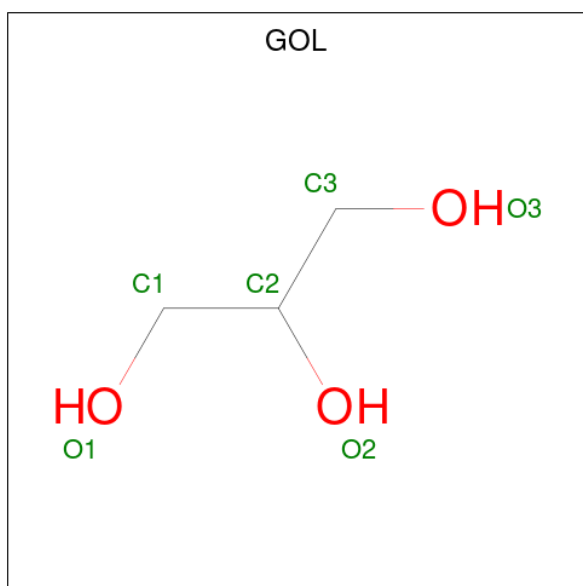
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Zn	0	0
			1	1		

- Molecule 6 is 2-AMINO-ETHANETHIOL (three-letter code: DHL) (formula: C₂H₇NS).



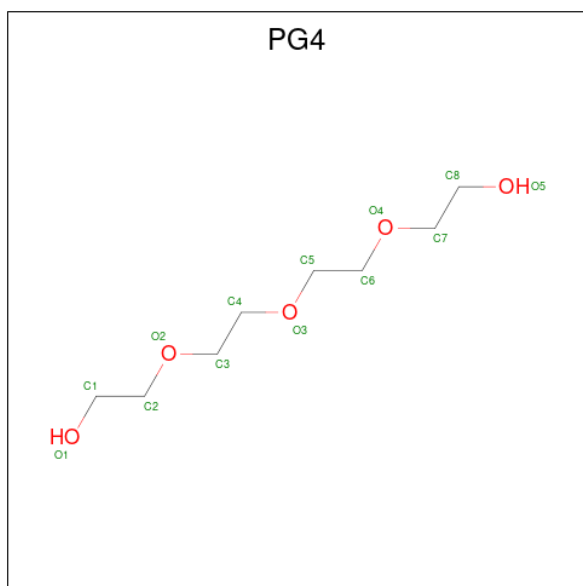
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	S		
6	A	1	4	2	1	1	0	0

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



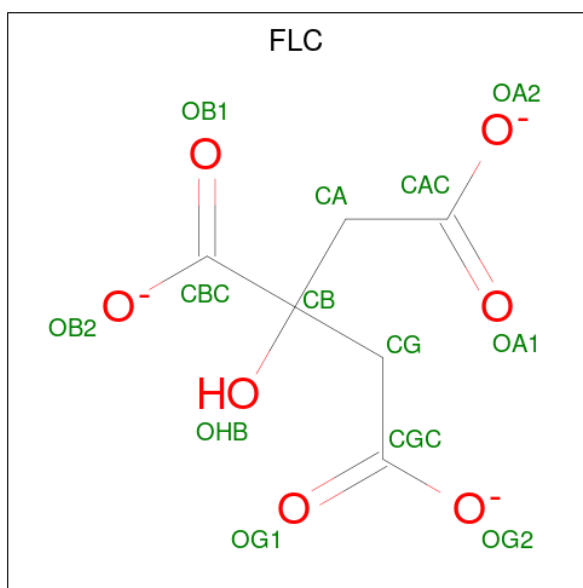
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 8 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			13	8	5		

- Molecule 9 is CITRATE ANION (three-letter code: FLC) (formula: $C_6H_5O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	C	O	0	0
			13	6	7		

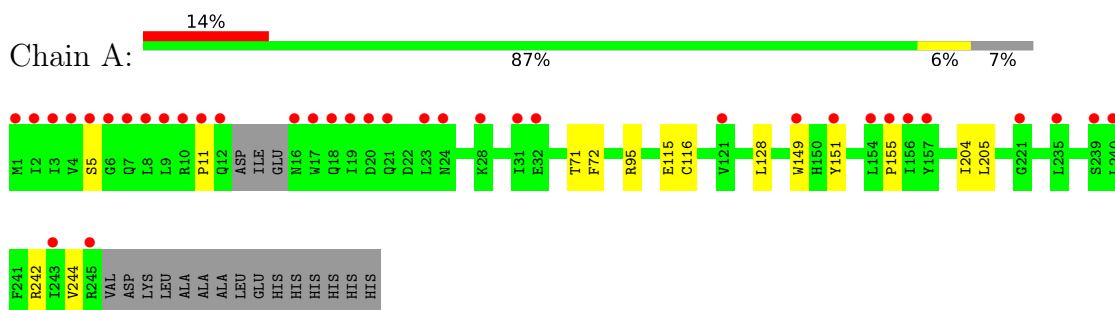
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	113	Total	O	0	3
			116	116		
10	B	91	Total	O	0	1
			92	92		

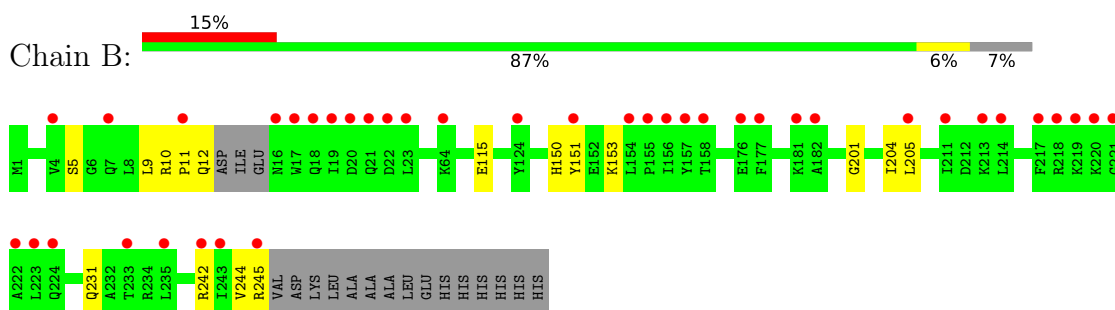
3 Residue-property plots [i](#)

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: Protein-glutamine gamma-glutamyltransferase



- Molecule 1: Protein-glutamine gamma-glutamyltransferase



4 Data and refinement statistics i

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	98.35Å 98.35Å 65.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.57 – 1.86 39.39 – 1.86	Depositor EDS
% Data completeness (in resolution range)	84.8 (36.57-1.86) 84.8 (39.39-1.86)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.75 (at 1.87Å)	Xtrriage
Refinement program	BUSTER-TNT 2.10, PHENIX 1.8.4-dev1565	Depositor
R, R_{free}	0.167 , 0.213 0.177 , 0.221	Depositor DCC
R_{free} test set	2237 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	30.8	Xtrriage
Anisotropy	0.842	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 60.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.032 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4269	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.95% 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: DHL, EDO, FLC, GOL, PGE, PG4, SO4, ZN

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	0.58	0/2044	0.63	0/2764
1	B	0.54	0/2048	0.62	0/2770
All	All	0.56	0/4092	0.62	0/5534

There are no bond length outliers.

There are no bond angle outliers.

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	1993	0	1997	8	0
1	B	1997	0	1996	8	0
2	A	20	0	28	1	0
3	A	4	0	6	2	0
4	A	5	0	0	0	0
4	B	5	0	0	1	0
5	A	1	0	0	0	0
6	A	4	0	6	2	0
7	A	6	0	8	0	0
8	A	13	0	18	0	0
9	B	13	0	5	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	A	116	0	0	0	0
10	B	92	0	0	1	0
All	All	4269	0	4064	14	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 2.

All (14) 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:116[B]:CYS:SG	6:A:2006:DHL:N	2.40	0.94
1:B:115[B]:GLU:OE1	10:B:7173[B]:HOH:O	1.98	0.81
1:B:10:ARG:O	1:B:12:GLN:N	2.44	0.51
1:A:116[B]:CYS:CB	6:A:2006:DHL:HN1	2.28	0.47
1:A:151:TYR:O	1:B:150:HIS:HA	2.15	0.46
1:A:72:PHE:H	3:A:2003:EDO:C1	2.30	0.45
1:B:201:GLY:N	4:B:7002:SO4:O3	2.42	0.43
1:A:242[B]:ARG:NH2	1:B:231:GLN:OE1	2.43	0.43
1:B:242:ARG:HA	1:B:242:ARG:HD2	1.89	0.43
1:A:71:THR:HB	3:A:2003:EDO:H11	2.01	0.42
1:A:149:TRP:CE2	2:A:2001:PGE:H5	2.55	0.42
1:A:204:ILE:C	1:A:205:LEU:HD12	2.41	0.41
1:B:204:ILE:C	1:B:205:LEU:HD12	2.41	0.40
1:B:244:VAL:O	1:B:245:ARG:HB3	2.22	0.40

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	240/260 (92%)	226 (94%)	12 (5%)	2 (1%)	19 7

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	241/260 (93%)	228 (95%)	10 (4%)	3 (1%)	13	3
All	All	481/520 (92%)	454 (94%)	22 (5%)	5 (1%)	14	5

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	11	PRO
1	B	5	SER
1	A	11	PRO
1	B	153	LYS
1	A	155	PRO

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	211/224 (94%)	206 (98%)	5 (2%)	49	33
1	B	212/224 (95%)	210 (99%)	2 (1%)	78	72
All	All	423/448 (94%)	416 (98%)	7 (2%)	60	47

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	SER
1	A	95	ARG
1	A	115	GLU
1	A	128	LEU
1	A	244	VAL
1	B	9	LEU
1	B	151	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

Of 10 ligands modelled in this entry, 1 is monoatomic - leaving 9 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	$\# Z > 2$
6	DHL	A	2006	1	2,3,3	0.48	0	1,2,2	0.31	0
4	SO4	B	7002	-	4,4,4	0.21	0	6,6,6	0.26	0
8	PG4	A	2008	-	12,12,12	0.40	0	11,11,11	0.62	0
4	SO4	A	2004	-	4,4,4	0.20	0	6,6,6	0.09	0
9	FLC	B	7001	-	12,12,12	1.00	0	17,17,17	1.72	2 (11%)
2	PGE	A	2002	-	9,9,9	0.46	0	8,8,8	0.41	0
7	GOL	A	2007	-	5,5,5	0.34	0	5,5,5	0.55	0
3	EDO	A	2003	-	3,3,3	0.64	0	2,2,2	0.06	0
2	PGE	A	2001	-	9,9,9	0.39	0	8,8,8	0.52	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
6	DHL	A	2006	1	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	PG4	A	2008	-	-	4/10/10/10	-
9	FLC	B	7001	-	-	5/16/16/16	-
2	PGE	A	2002	-	-	5/7/7/7	-
7	GOL	A	2007	-	-	2/4/4/4	-
3	EDO	A	2003	-	-	1/1/1/1	-
2	PGE	A	2001	-	-	2/7/7/7	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	B	7001	FLC	OB2-CBC-CB	4.21	120.36	113.05
9	B	7001	FLC	OB1-CBC-CB	-4.04	116.53	122.25

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	B	7001	FLC	CAC-CA-CB-CBC
9	B	7001	FLC	CAC-CA-CB-CG
9	B	7001	FLC	CAC-CA-CB-OHB
2	A	2002	PGE	O2-C3-C4-O3
8	A	2008	PG4	O2-C3-C4-O3
9	B	7001	FLC	CB-CG-CGC-OG1
9	B	7001	FLC	CB-CG-CGC-OG2
7	A	2007	GOL	O1-C1-C2-O2
2	A	2002	PGE	C3-C4-O3-C5
2	A	2001	PGE	O3-C5-C6-O4
2	A	2001	PGE	C6-C5-O3-C4
8	A	2008	PG4	C4-C3-O2-C2
2	A	2002	PGE	C1-C2-O2-C3
3	A	2003	EDO	O1-C1-C2-O2
8	A	2008	PG4	C6-C5-O3-C4
2	A	2002	PGE	O1-C1-C2-O2
7	A	2007	GOL	O1-C1-C2-C3
2	A	2002	PGE	O3-C5-C6-O4
8	A	2008	PG4	O4-C7-C8-O5

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	2006	DHL	2	0
4	B	7002	SO4	1	0
3	A	2003	EDO	2	0
2	A	2001	PGE	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	242/260 (93%)	0.63	36 (14%) 2 2	24, 42, 100, 145	0
1	B	242/260 (93%)	0.74	40 (16%) 1 1	25, 47, 95, 113	0
All	All	484/520 (93%)	0.68	76 (15%) 2 2	24, 44, 100, 145	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	223	LEU	11.1
1	A	8	LEU	8.3
1	A	9	LEU	6.6
1	B	222	ALA	5.6
1	A	17	TRP	5.4
1	B	18	GLN	5.3
1	B	220	LYS	5.2
1	A	11	PRO	5.1
1	B	182	ALA	4.9
1	A	10	ARG	4.7
1	B	224	GLN	4.6
1	A	156	ILE	4.5
1	B	16	ASN	4.5
1	B	221	GLY	4.5
1	A	21	GLN	4.5
1	B	7	GLN	4.5
1	B	17	TRP	4.4
1	A	245	ARG	4.3
1	A	1	MET	4.3
1	B	156	ILE	3.9
1	B	154	LEU	3.9
1	A	7	GLN	3.8
1	B	218	ARG	3.7
1	A	3	ILE	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	151	TYR	3.6
1	A	4	VAL	3.6
1	A	2	ILE	3.5
1	A	18	GLN	3.5
1	A	16	ASN	3.4
1	A	154	LEU	3.4
1	B	151	TYR	3.4
1	B	217	PHE	3.3
1	A	31	ILE	3.2
1	B	243	ILE	3.2
1	A	235	LEU	3.1
1	B	245	ARG	3.1
1	B	219	LYS	3.1
1	A	12	GLN	3.1
1	A	23	LEU	2.9
1	A	6	GLY	2.9
1	B	214	LEU	2.9
1	B	211	ILE	2.9
1	B	177	PHE	2.9
1	A	5	SER	2.8
1	B	176	GLU	2.8
1	B	23	LEU	2.8
1	A	19	ILE	2.8
1	B	213	LYS	2.7
1	A	243	ILE	2.7
1	B	235	LEU	2.7
1	A	24	ASN	2.6
1	B	157	TYR	2.6
1	A	32	GLU	2.5
1	B	22	ASP	2.5
1	B	20	ASP	2.5
1	B	181	LYS	2.4
1	A	121	VAL	2.4
1	A	240	LEU	2.4
1	B	205	LEU	2.3
1	A	221	GLY	2.3
1	B	4	VAL	2.3
1	B	64	LYS	2.3
1	B	19	ILE	2.3
1	A	149	TRP	2.2
1	A	155	PRO	2.2
1	B	21	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	28	LYS	2.2
1	B	124	TYR	2.2
1	B	233	THR	2.1
1	B	11	PRO	2.1
1	B	242	ARG	2.1
1	A	239	SER	2.1
1	A	157	TYR	2.1
1	B	158	THR	2.1
1	B	155	PRO	2.0
1	A	20	ASP	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
7	GOL	A	2007	6/6	0.69	0.18	67,68,71,79	0
3	EDO	A	2003	4/4	0.76	0.16	34,59,67,76	0
9	FLC	B	7001	13/13	0.81	0.21	59,80,87,87	0
6	DHL	A	2006	4/4	0.85	0.16	40,45,50,76	4
8	PG4	A	2008	13/13	0.89	0.24	42,52,71,74	0
2	PGE	A	2002	10/10	0.92	0.10	40,52,71,72	0
4	SO4	B	7002	5/5	0.94	0.17	77,78,86,94	0
5	ZN	A	2005	1/1	0.94	0.08	80,80,80,80	0
4	SO4	A	2004	5/5	0.95	0.26	84,94,101,107	0
2	PGE	A	2001	10/10	0.96	0.17	43,51,61,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.