



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:56 AM UTC

PDB ID : 6ZOE / pdb_00006zoe
Title : AcrB-F563A symmetric T protomer
Authors : Tam, H.K.; Foong, W.E.; Pos, K.M.
Deposited on : 2020-07-07
Resolution : 2.85 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 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

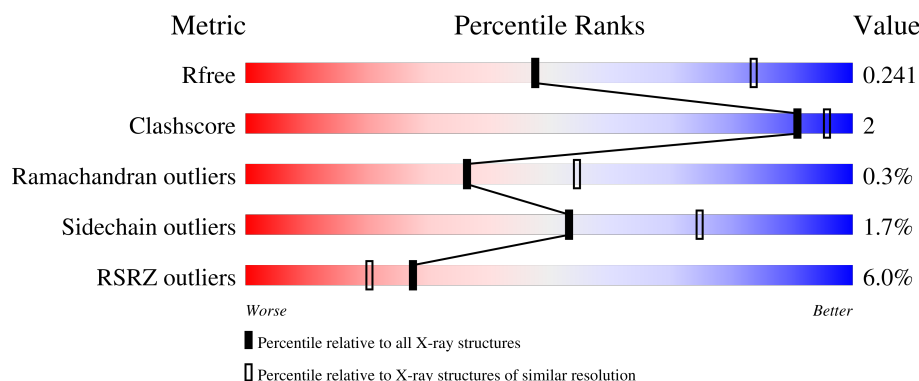
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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}	180053	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	1057	6% 92% 6%
2	B	169	4% 91% 8%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 9224 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 Multidrug efflux pump subunit AcrB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1040	Total	C	N	O	S	0	2	0
			7918	5090	1313	1471	44			

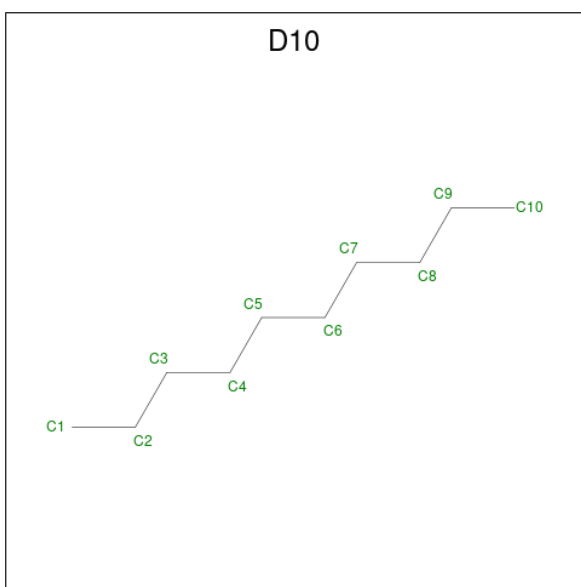
There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	563	ALA	PHE	engineered mutation	UNP P31224
A	1050	LEU	-	expression tag	UNP P31224
A	1051	GLU	-	expression tag	UNP P31224
A	1052	HIS	-	expression tag	UNP P31224
A	1053	HIS	-	expression tag	UNP P31224
A	1054	HIS	-	expression tag	UNP P31224
A	1055	HIS	-	expression tag	UNP P31224
A	1056	HIS	-	expression tag	UNP P31224
A	1057	HIS	-	expression tag	UNP P31224

- Molecule 2 is a protein called DARPIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	156	Total	C	N	O	S	0	0	0
			1177	741	206	229	1			

- Molecule 3 is DECANE (CCD ID: D10) (formula: C₁₀H₂₂).



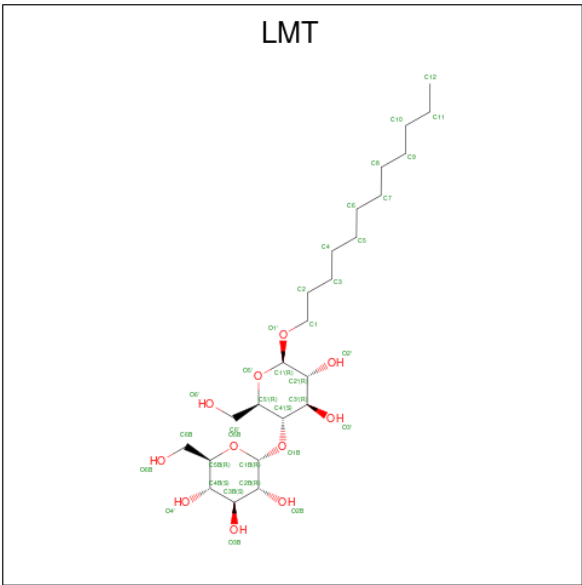
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C	0	0
			10	10		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



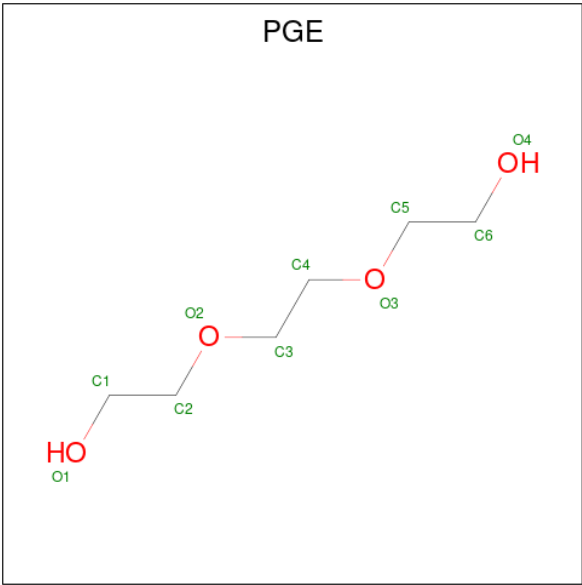
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is DODECYL-BETA-D-MALTOSE (CCD ID: LMT) (formula: $C_{24}H_{46}O_{11}$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			35	24	11		

- Molecule 6 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



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

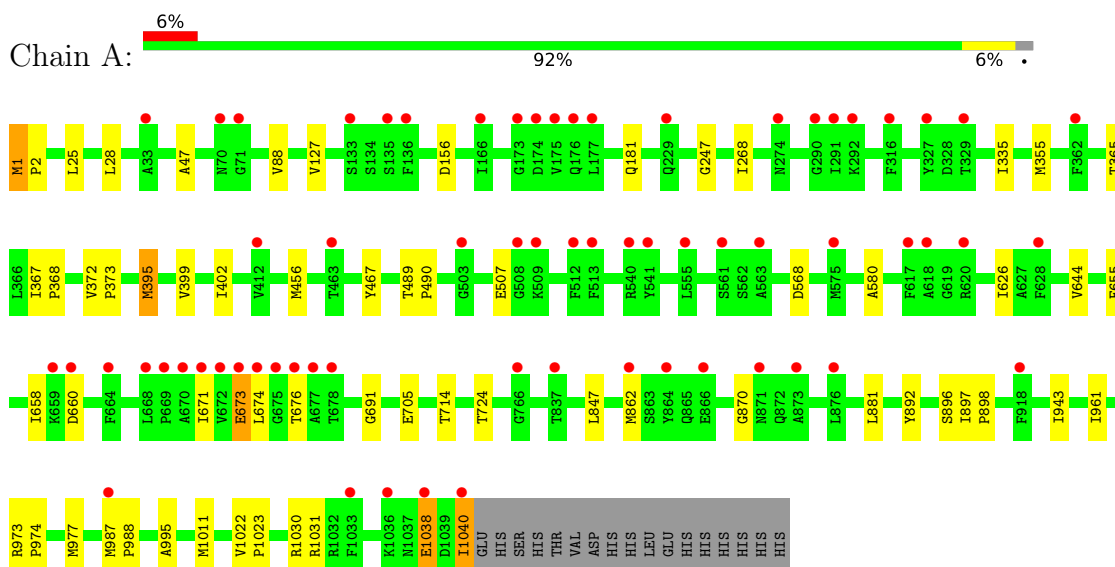
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	41	Total 43	O 43	0	2
7	B	13	Total 13	O 13	0	0

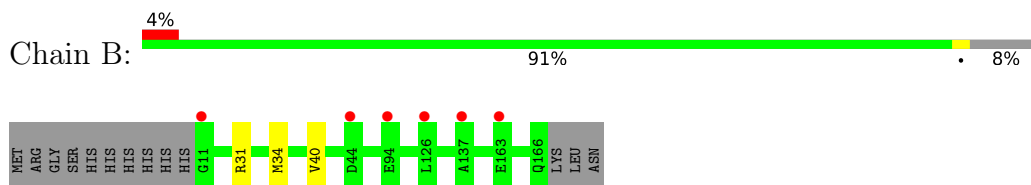
3 Residue-property plots [i](#)

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: Multidrug efflux pump subunit AcrB



- Molecule 2: DARPIN



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, α , β , γ	226.61Å 226.61Å 226.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.26 – 2.85 46.26 – 2.85	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.26-2.85) 100.0 (46.26-2.85)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.17 (at 2.86Å)	Xtriage
Refinement program	REFMAC 5.8.0253	Depositor
R, R_{free}	0.223 , 0.265 (Not available) , 0.241	Depositor DCC
R_{free} test set	2320 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å ²)	70.5	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 39.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.004 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9224	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.12% 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: LMT, EDO, D10, PGE

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	1.03	0/8073	1.55	3/10960 (0.0%)
2	B	1.03	0/1196	1.59	0/1626
All	All	1.03	0/9269	1.56	3/12586 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	896	SER	CA-C-N	7.67	125.07	120.24
1	A	896	SER	C-N-CA	7.67	125.07	120.24
1	A	691	GLY	CA-C-O	-5.66	118.54	122.22

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	7918	0	8081	30	0
2	B	1177	0	1159	1	0
3	A	10	0	22	0	0
4	A	4	0	6	0	0
4	B	4	0	6	0	0
5	A	35	0	46	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	20	0	28	0	0
7	A	43	0	0	0	0
7	B	13	0	0	0	0
All	All	9224	0	9348	31	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 (31) 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:580:ALA:HB1	1:A:724:THR:HG22	1.87	0.57
1:A:395:MET:HA	1:A:395:MET:HE2	1.90	0.53
1:A:1:MET:HB3	1:A:2:PRO:HD3	1.90	0.52
1:A:355:MET:HE2	1:A:368:PRO:HG2	1.92	0.51
1:A:335:ILE:HG21	1:A:995:ALA:HB3	1.93	0.51
1:A:372:VAL:HB	1:A:373:PRO:HD3	1.94	0.51
1:A:973:ARG:HG2	1:A:977:MET:HE2	1.93	0.49
1:A:1022:VAL:N	1:A:1023:PRO:HD2	2.29	0.48
1:A:705:GLU:HB3	1:A:847:LEU:HD22	1.96	0.47
1:A:367:ILE:HB	1:A:368:PRO:HD3	1.97	0.47
1:A:399:VAL:O	1:A:402:ILE:HG22	2.15	0.47
1:A:456:MET:HG2	1:A:467:TYR:HB3	1.98	0.46
1:A:1011:MET:HA	1:A:1011:MET:HE2	1.96	0.46
1:A:568:ASP:CG	1:A:644:VAL:HG23	2.41	0.46
1:A:671:ILE:HG22	1:A:673:GLU:HG3	1.97	0.46
1:A:489:THR:N	1:A:490:PRO:CD	2.79	0.46
1:A:961:ILE:HD11	1:A:1031:ARG:NH2	2.33	0.44
1:A:489:THR:OG1	1:A:490:PRO:HD3	2.18	0.43
1:A:47:ALA:HB2	1:A:127:VAL:HG13	2.00	0.42
2:B:34:MET:HE1	2:B:40:VAL:HG12	2.00	0.42
1:A:973:ARG:N	1:A:974:PRO:HD2	2.34	0.42
1:A:156:ASP:HA	1:A:181:GLN:HA	2.01	0.42
1:A:987:MET:N	1:A:988:PRO:CD	2.83	0.42
1:A:247:GLY:HA2	1:A:268:ILE:CD1	2.50	0.42
1:A:655:PHE:HA	1:A:658:ILE:HG12	2.02	0.41
1:A:892:TYR:OH	1:A:943:ILE:HA	2.20	0.41
1:A:897:ILE:N	1:A:898:PRO:CD	2.83	0.41
1:A:355:MET:HE2	1:A:368:PRO:CG	2.50	0.41
1:A:580:ALA:CB	1:A:724:THR:HG22	2.50	0.41
1:A:674:LEU:HD12	1:A:674:LEU:HA	1.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1038:GLU:HG3	1:A:1040:ILE:H	1.86	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	1040/1057 (98%)	1000 (96%)	37 (4%)	3 (0%)	36	54
2	B	154/169 (91%)	152 (99%)	2 (1%)	0	100	100
All	All	1194/1226 (97%)	1152 (96%)	39 (3%)	3 (0%)	36	54

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	676	THR
1	A	660	ASP
1	A	870	GLY

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	847/862 (98%)	832 (98%)	15 (2%)	51	73
2	B	120/132 (91%)	119 (99%)	1 (1%)	73	86

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	967/994 (97%)	951 (98%)	16 (2%)	53 75

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

Mol	Chain	Res	Type
1	A	1	MET
1	A	25	LEU
1	A	28	LEU
1	A	88	VAL
1	A	365	THR
1	A	395	MET
1	A	507	GLU
1	A	626	ILE
1	A	673	GLU
1	A	714	THR
1	A	862	MET
1	A	881	LEU
1	A	1030	ARG
1	A	1038	GLU
1	A	1040	ILE
2	B	31	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	GLN
1	A	125	GLN
1	A	469	GLN
1	A	1000	GLN
1	A	1037	ASN
2	B	125	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry

6 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	$\# Z > 2$
5	LMT	A	1103	-	36,36,36	0.45	0	47,47,47	0.64	0
4	EDO	B	301	-	3,3,3	0.07	0	2,2,2	0.06	0
6	PGE	A	1104	-	9,9,9	0.20	0	8,8,8	0.09	0
6	PGE	A	1105	-	9,9,9	0.17	0	8,8,8	0.11	0
3	D10	A	1101	-	9,9,9	0.11	0	8,8,8	0.07	0
4	EDO	A	1102	-	3,3,3	0.06	0	2,2,2	0.10	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
5	LMT	A	1103	-	-	9/21/61/61	0/2/2/2
4	EDO	B	301	-	-	1/1/1/1	-
6	PGE	A	1104	-	-	3/7/7/7	-
6	PGE	A	1105	-	-	3/7/7/7	-
3	D10	A	1101	-	-	3/7/7/7	-
4	EDO	A	1102	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

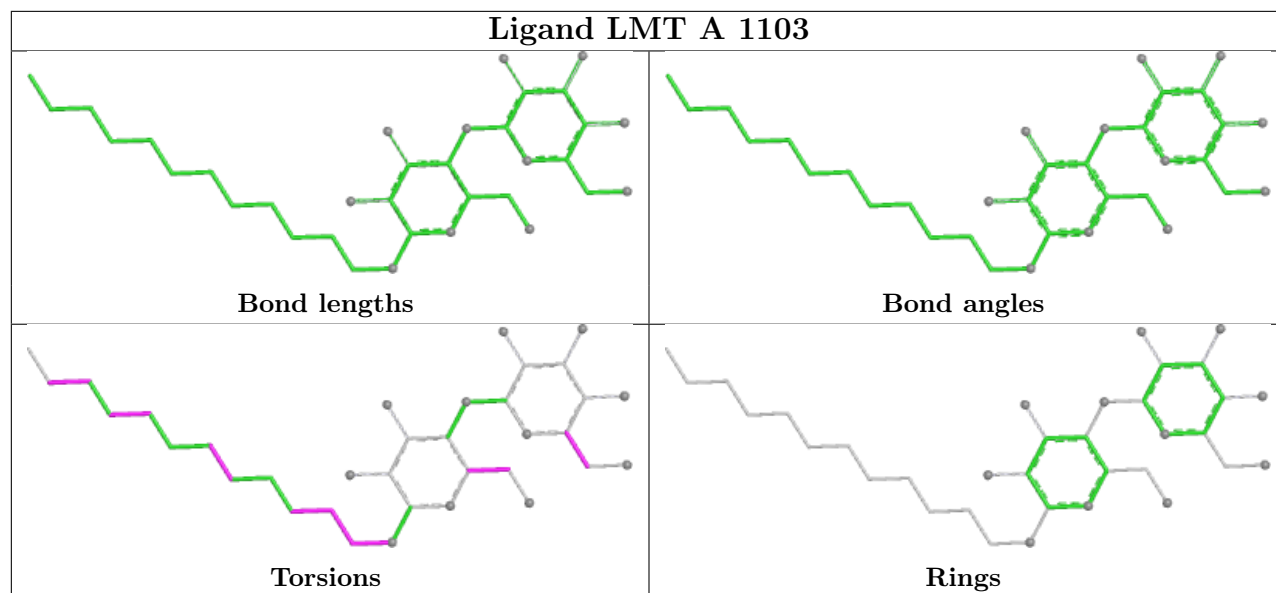
All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1103	LMT	O5B-C5B-C6B-O6B
5	A	1103	LMT	O1'-C1-C2-C3
5	A	1103	LMT	C7-C8-C9-C10
6	A	1104	PGE	O1-C1-C2-O2
5	A	1103	LMT	C1-C2-C3-C4
4	A	1102	EDO	O1-C1-C2-O2
5	A	1103	LMT	C4-C5-C6-C7
6	A	1105	PGE	O1-C1-C2-O2
5	A	1103	LMT	C9-C10-C11-C12
6	A	1104	PGE	C6-C5-O3-C4
6	A	1104	PGE	O2-C3-C4-O3
3	A	1101	D10	C1-C2-C3-C4
5	A	1103	LMT	O5'-C5'-C6'-O6'
5	A	1103	LMT	C2-C1-O1'-C1'
6	A	1105	PGE	O2-C3-C4-O3
3	A	1101	D10	C3-C4-C5-C6
4	B	301	EDO	O1-C1-C2-O2
5	A	1103	LMT	C4B-C5B-C6B-O6B
6	A	1105	PGE	C4-C3-O2-C2
3	A	1101	D10	C5-C6-C7-C8

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	1040/1057 (98%)	0.51	66 (6%) 26 19	47, 83, 109, 125	2 (0%)
2	B	156/169 (92%)	0.30	6 (3%) 44 35	54, 74, 123, 140	0
All	All	1196/1226 (97%)	0.49	72 (6%) 27 20	47, 82, 110, 140	2 (0%)

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	174	ASP	6.7
1	A	676	THR	6.7
1	A	135	SER	6.5
1	A	866	GLU	5.9
1	A	292	LYS	5.2
1	A	136	PHE	5.0
1	A	672	VAL	4.9
1	A	675	GLY	4.8
1	A	175	VAL	4.5
1	A	1040	ILE	4.4
1	A	508	GLY	4.3
1	A	987	MET	4.2
1	A	541	TYR	3.9
1	A	677	ALA	3.6
1	A	561	SER	3.6
1	A	1033	PHE	3.6
1	A	673	GLU	3.5
1	A	664	PHE	3.4
1	A	660	ASP	3.3
1	A	166	ILE	3.3
1	A	327	TYR	3.2
1	A	659	LYS	3.1
1	A	1036	LYS	3.1
1	A	316	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	670	ALA	2.9
1	A	671	ILE	2.9
1	A	290	GLY	2.9
1	A	133	SER	2.9
1	A	291	ILE	2.8
1	A	1038	GLU	2.7
1	A	513	PHE	2.7
1	A	620	ARG	2.7
1	A	563	ALA	2.6
1	A	668	LEU	2.6
2	B	94	GLU	2.6
1	A	617	PHE	2.6
1	A	628	PHE	2.6
1	A	509	LYS	2.5
1	A	176	GLN	2.5
1	A	503	GLY	2.5
1	A	412	VAL	2.5
1	A	540	ARG	2.5
1	A	618	ALA	2.5
1	A	837	THR	2.4
1	A	669	PRO	2.4
1	A	876	LEU	2.4
1	A	674	LEU	2.4
1	A	229	GLN	2.3
1	A	862	MET	2.3
1	A	463	THR	2.3
1	A	864	TYR	2.3
1	A	70	ASN	2.2
1	A	678	THR	2.2
1	A	766	GLY	2.2
1	A	871	ASN	2.2
1	A	512	PHE	2.2
1	A	33	ALA	2.2
2	B	137	ALA	2.2
2	B	126	LEU	2.1
1	A	274	ASN	2.1
2	B	44	ASP	2.1
2	B	11	GLY	2.1
1	A	918	PHE	2.1
1	A	575	MET	2.1
1	A	71	GLY	2.1
1	A	362	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	873	ALA	2.1
2	B	163	GLU	2.1
1	A	555	LEU	2.0
1	A	177	LEU	2.0
1	A	173	GLY	2.0
1	A	329	THR	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

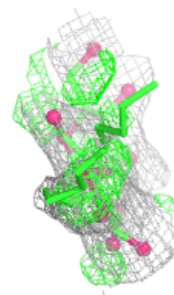
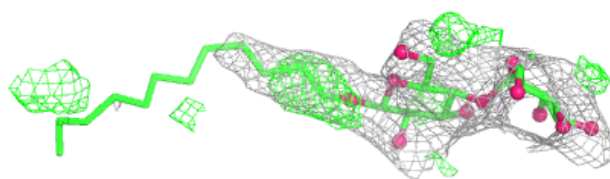
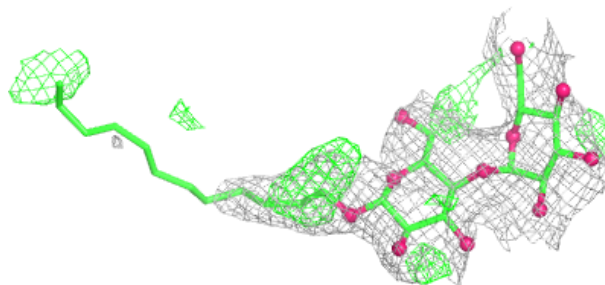
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	EDO	B	301	4/4	0.79	0.22	100,101,103,105	0
6	PGE	A	1104	10/10	0.82	0.17	90,94,96,96	0
3	D10	A	1101	10/10	0.83	0.32	100,103,106,106	0
4	EDO	A	1102	4/4	0.84	0.32	88,89,90,91	0
5	LMT	A	1103	35/35	0.86	0.26	88,91,93,94	35
6	PGE	A	1105	10/10	0.93	0.15	89,92,96,100	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around LMT A 1103:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.