



wwPDB X-ray Structure Validation Summary Report ⓘ

May 15, 2020 – 01:10 pm BST

PDB ID : 6EER
Title : Structure of glycine-bound GoxA from *Pseudoalteromonas luteoviolacea*
Authors : Yukl, E.T.; Avalos, D.
Deposited on : 2018-08-15
Resolution : 1.82 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

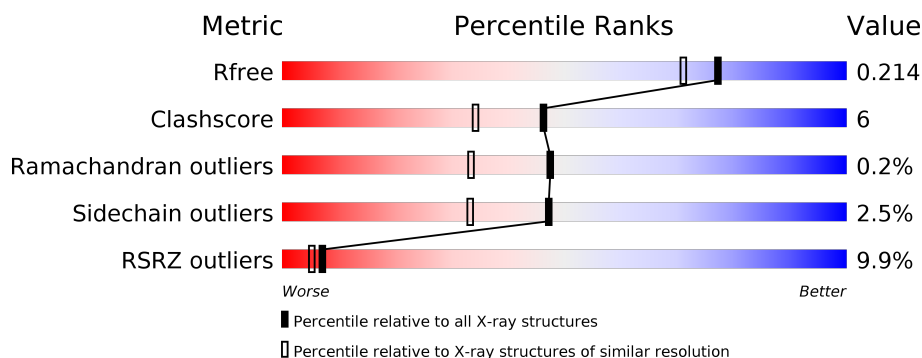
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.82 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	7484 (1.84-1.80)
Clashscore	141614	8401 (1.84-1.80)
Ramachandran outliers	138981	8290 (1.84-1.80)
Sidechain outliers	138945	8290 (1.84-1.80)
RSRZ outliers	127900	7371 (1.84-1.80)

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	816	<div> <div>8%</div> <div>83%</div> <div>10%</div> <div>6%</div> </div>
1	B	816	<div> <div>11%</div> <div>83%</div> <div>13%</div> <div>•</div> </div>
1	C	816	<div> <div>8%</div> <div>82%</div> <div>12%</div> <div>5%</div> </div>
1	D	816	<div> <div>11%</div> <div>82%</div> <div>13%</div> <div>• •</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	PEG	C	903	-	-	X	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 27935 atoms, of which 30 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 GoxA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	768	Total	C	N	O	S	0	14	0
			6161	3891	1049	1201	20			
1	B	786	Total	C	N	O	S	0	10	0
			6298	3976	1073	1228	21			
1	C	777	Total	C	N	O	S	0	7	0
			6202	3921	1057	1204	20			
1	D	784	Total	C	N	O	S	0	5	0
			6241	3941	1062	1218	20			

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		
2	D	1	Total	Mg	0	0
			1	1		
2	C	1	Total	Mg	0	0
			1	1		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

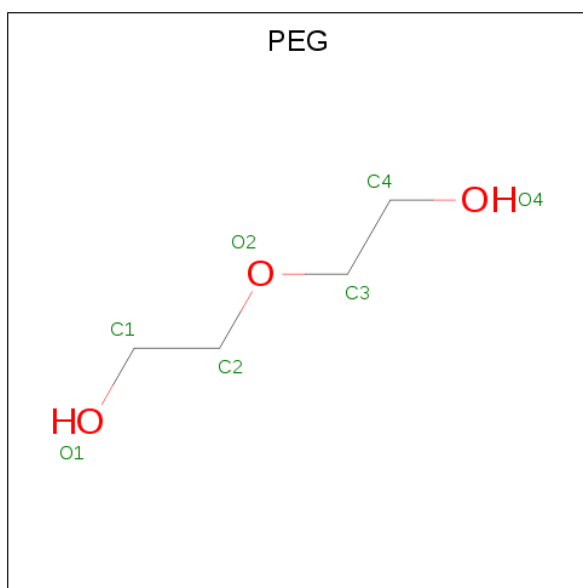
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Na	0	0
			1	1		
3	A	1	Total	Na	0	0
			1	1		
3	D	1	Total	Na	0	0
			1	1		
3	C	1	Total	Na	0	0
			1	1		

- Molecule 4 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			24	6	14	4		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	C	1	Total	C	H	O	0	0
			17	4	10	3		

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	D	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	776	Total	O	0	5
			781	781		
7	B	699	Total	O	0	5
			704	704		
7	C	772	Total	O	0	3
			775	775		
7	D	705	Total	O	0	9
			714	714		

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.

Chain A:

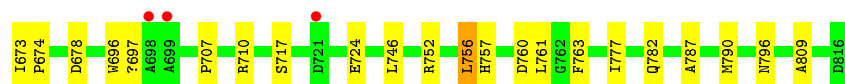
8% 10% 6%

83%

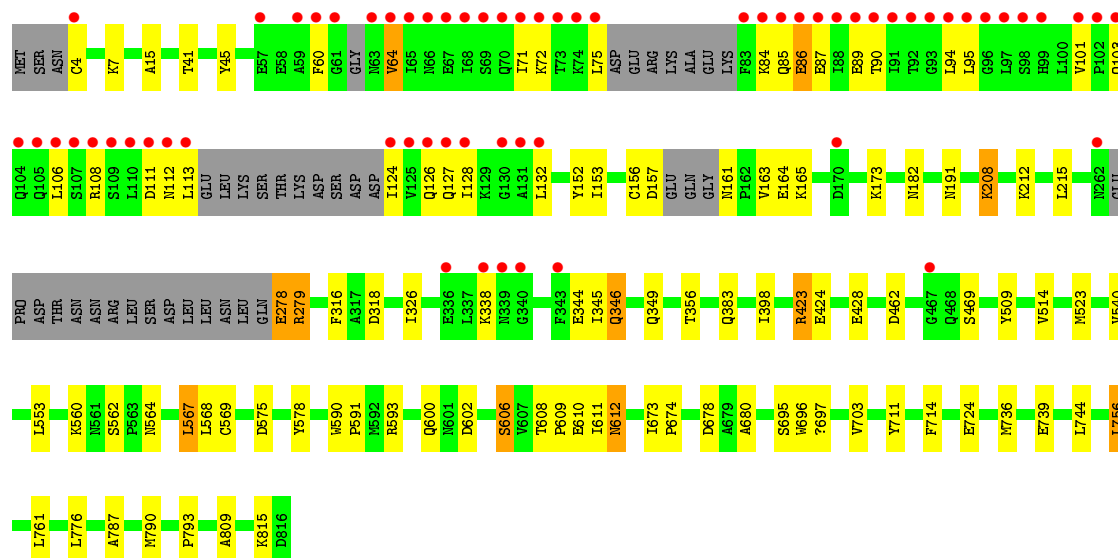
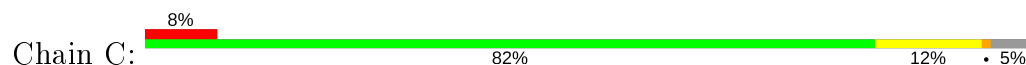
MET SER ASN C4 Q5 A15 K24 Y45 Q49 L56 E57 E58 A59 F60 G61 G62 N63 V64 I65 M66 E67 I68 S69 Q70 I71 K72 THR LVS LEU ASP GLU ARG LVS ALA GIU LVS PHE LYS GLN GIU I88 E89 T90 I91 T92 G93 I94 L95 G96 L97 S98 H99 I100 V101

Chain B:

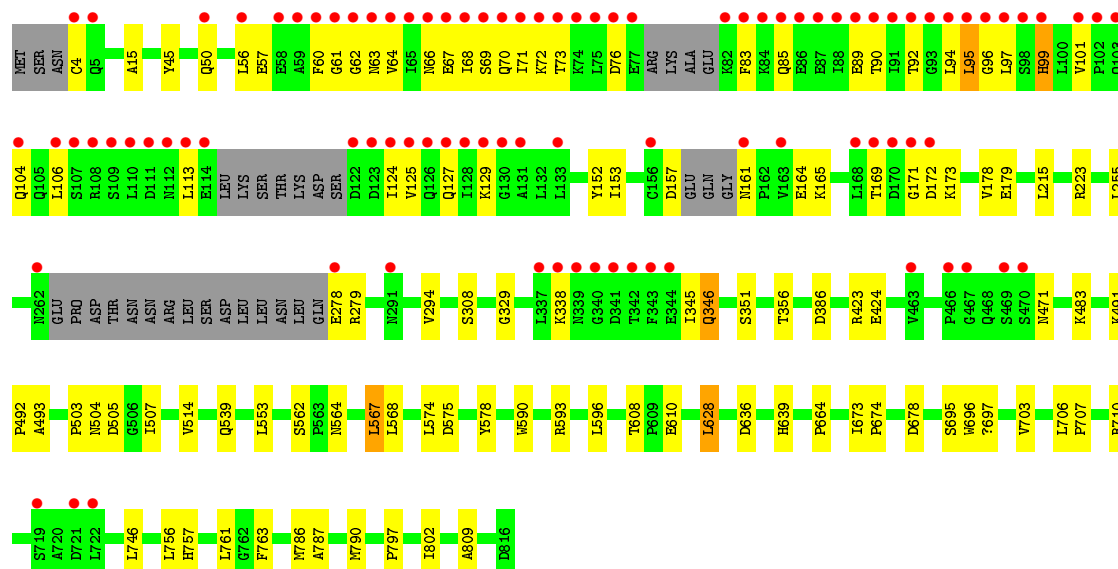
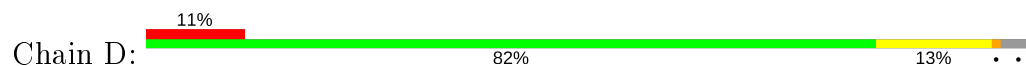
Category	Percentage
Green	11%
Yellow	83%
Red	13%



• Molecule 1: GoxA



• Molecule 1: GoxA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	110.24Å 93.46Å 187.90Å 90.00° 95.02° 90.00°	Depositor
Resolution (Å)	48.30 – 1.82 48.30 – 1.82	Depositor EDS
% Data completeness (in resolution range)	99.4 (48.30-1.82) 99.4 (48.30-1.82)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 1.82Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.175 , 0.213 0.177 , 0.214	Depositor DCC
R_{free} test set	17055 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.476	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 51.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	27935	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.37% 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: MG, PGE, TNQ, NA, EDO, PEG

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	0.38	0/6287	0.56	0/8563
1	B	0.35	0/6425	0.54	0/8739
1	C	0.38	0/6328	0.57	0/8612
1	D	0.36	0/6368	0.55	0/8667
All	All	0.37	0/25408	0.56	0/34581

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	6161	0	5903	59	0
1	B	6298	0	6048	82	0
1	C	6202	0	5957	83	0
1	D	6241	0	5984	83	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	10	14	14	0	0
5	C	7	10	10	8	0
6	D	4	6	6	3	0
7	A	781	0	0	11	0
7	B	704	0	0	12	0
7	C	775	0	0	13	0
7	D	714	0	0	8	0
All	All	27905	30	23922	308	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.

The worst 5 of 308 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:D:67:GLU:HG2	1:D:94:LEU:HD11	1.28	1.15
1:B:179:GLU:HB2	1:B:628:LEU:HD22	1.54	0.88
1:D:179:GLU:HB2	1:D:628:LEU:HD22	1.56	0.86
1:D:83:PHE:HZ	1:D:124:ILE:HD11	1.41	0.84
1:C:191:ASN:HA	5:C:903:PEG:H31	1.61	0.83

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	771/816 (94%)	747 (97%)	23 (3%)	1 (0%)	51 37
1	B	785/816 (96%)	764 (97%)	19 (2%)	2 (0%)	41 27

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	C	771/816 (94%)	750 (97%)	20 (3%)	1 (0%)	51 37
1	D	778/816 (95%)	754 (97%)	22 (3%)	2 (0%)	41 27
All	All	3105/3264 (95%)	3015 (97%)	84 (3%)	6 (0%)	47 33

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62	GLY
1	B	85	GLN
1	B	279	ARG
1	D	96	GLY
1	D	279	ARG

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	679/711 (96%)	664 (98%)	15 (2%)	52 39
1	B	693/711 (98%)	682 (98%)	11 (2%)	62 53
1	C	682/711 (96%)	658 (96%)	24 (4%)	36 20
1	D	686/711 (96%)	664 (97%)	22 (3%)	39 24
All	All	2740/2844 (96%)	2668 (97%)	72 (3%)	47 32

5 of 72 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	278	GLU
1	C	567	LEU
1	D	578	TYR
1	C	338	LYS
1	C	423[B]	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	103	GLN
1	C	127	GLN
1	C	346	GLN
1	D	99	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues 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$
1	TNQ	D	697	1	15,21,22	1.94	3 (20%)	13,29,31	2.80	5 (38%)
1	TNQ	C	697	1	15,21,22	1.87	4 (26%)	13,29,31	2.62	3 (23%)
1	TNQ	B	697	1	15,21,22	1.99	3 (20%)	13,29,31	2.58	4 (30%)
1	TNQ	A	697	1	15,21,22	1.94	4 (26%)	13,29,31	2.54	3 (23%)

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
1	TNQ	D	697	1	-	2/7/11/13	0/2/2/2
1	TNQ	C	697	1	-	2/7/11/13	0/2/2/2
1	TNQ	B	697	1	-	2/7/11/13	0/2/2/2
1	TNQ	A	697	1	-	2/7/11/13	0/2/2/2

The worst 5 of 14 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	697	TNQ	C2-N1	-5.30	1.27	1.45
1	B	697	TNQ	C2-N1	-5.30	1.27	1.45
1	D	697	TNQ	C2-N1	-5.27	1.28	1.45
1	A	697	TNQ	C2-N1	-5.25	1.28	1.45
1	D	697	TNQ	CZ2-CE2	3.31	1.48	1.42

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	697	TNQ	C1-C2-N1	8.15	126.15	110.43
1	B	697	TNQ	C1-C2-N1	7.81	125.49	110.43
1	C	697	TNQ	C1-C2-N1	7.56	125.00	110.43
1	A	697	TNQ	C1-C2-N1	7.27	124.45	110.43
1	A	697	TNQ	C2-N1-CH2	-4.25	111.11	121.97

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	D	697	TNQ	CZ3-CH2-N1-C2
1	B	697	TNQ	CZ3-CH2-N1-C2
1	D	697	TNQ	CZ2-CH2-N1-C2
1	B	697	TNQ	CZ2-CH2-N1-C2
1	A	697	TNQ	CZ2-CH2-N1-C2

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	697	TNQ	2	0
1	C	697	TNQ	2	0
1	B	697	TNQ	2	0
1	A	697	TNQ	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 8 are monoatomic - leaving 3 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	$\# Z > 2$
4	PGE	A	903	-	9,9,9	0.40	0	8,8,8	0.52	0
5	PEG	C	903	-	6,6,6	0.42	0	5,5,5	0.46	0
6	EDO	D	903	-	3,3,3	0.54	0	2,2,2	0.20	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
4	PGE	A	903	-	-	3/7/7/7	-
5	PEG	C	903	-	-	1/4/4/4	-
6	EDO	D	903	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	903	PEG	O1-C1-C2-O2
4	A	903	PGE	O2-C3-C4-O3
4	A	903	PGE	O1-C1-C2-O2
4	A	903	PGE	C4-C3-O2-C2

There are no ring outliers.

2 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	903	PEG	8	0
6	D	903	EDO	3	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 [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, 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	767/816 (93%)	0.30	62 (8%) 12 9	12, 22, 71, 120	0
1	B	785/816 (96%)	0.40	89 (11%) 5 3	15, 26, 72, 117	0
1	C	776/816 (95%)	0.29	64 (8%) 11 9	13, 22, 72, 122	0
1	D	783/816 (95%)	0.47	93 (11%) 4 3	14, 26, 75, 125	0
All	All	3111/3264 (95%)	0.37	308 (9%) 7 5	12, 24, 73, 125	0

The worst 5 of 308 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	71	ILE	11.9
1	D	124	ILE	10.7
1	A	71	ILE	10.2
1	A	64	VAL	9.9
1	B	115	LEU	9.7

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	TNQ	B	697	20/21	0.93	0.12	13,17,26,30	0
1	TNQ	C	697	20/21	0.94	0.12	12,17,25,28	0
1	TNQ	A	697	20/21	0.95	0.10	11,16,25,25	0
1	TNQ	D	697	20/21	0.96	0.11	14,18,28,30	0

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(\AA^2)	Q<0.9
4	PGE	A	903	10/10	0.67	0.25	51,65,74,76	0
5	PEG	C	903	7/7	0.71	0.27	41,54,67,67	0
6	EDO	D	903	4/4	0.74	0.39	34,42,50,60	0
3	NA	B	902	1/1	0.96	0.14	45,45,45,45	0
3	NA	A	902	1/1	0.96	0.18	38,38,38,38	0
3	NA	C	902	1/1	0.96	0.23	41,41,41,41	0
2	MG	D	901	1/1	0.98	0.08	23,23,23,23	0
3	NA	D	902	1/1	0.98	0.07	37,37,37,37	0
2	MG	B	901	1/1	0.98	0.08	24,24,24,24	0
2	MG	C	901	1/1	0.99	0.12	18,18,18,18	0
2	MG	A	901	1/1	0.99	0.08	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.