wwPDB X-ray Structure Validation Summary Report (i)

Jun 24, 2024 – 07:35 AM EDT

PDB ID	:	6ESZ
Title	:	Crystal structure of PqsBC from Pseudomonas aeruginosa (crystal form 1)
Authors	:	Witzgall, F.; Blankenfeldt, W.
Deposited on	:	2017-10-25
Resolution	:	1.84 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 (i) 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 (1)) were used in the production of this report:

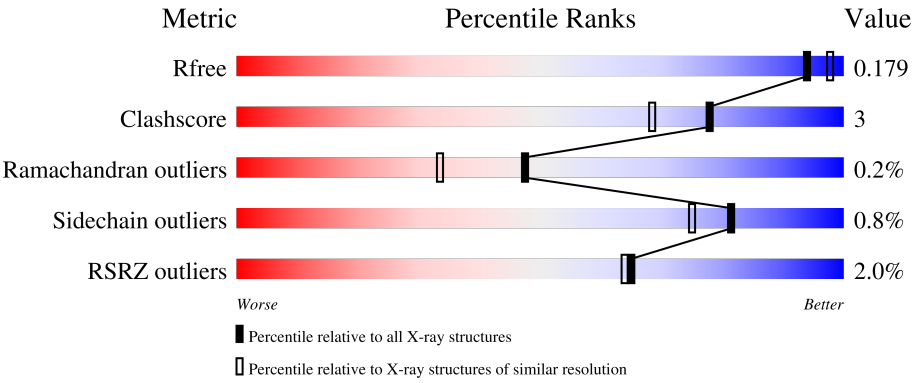
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.84 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	4003 (1.86-1.82)
Clashscore	141614	4233 (1.86-1.82)
Ramachandran outliers	138981	4185 (1.86-1.82)
Sidechain outliers	138945	4186 (1.86-1.82)
RSRZ outliers	127900	3957 (1.86-1.82)

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 <=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	365	2% 93% 5%
1	C	365	3% 90% 7%
2	B	283	% 94% 5%
2	D	283	% 96%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 21569 atoms, of which 10211 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 PqsC.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	355	Total	C	H	N	O	S	0	13	0
			5619	1788	2805	483	521	22			
1	C	354	Total	C	H	N	O	S	0	13	0
			5555	1773	2770	471	519	22			

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	MET	-	initiating methionine	UNP Q9I4X1
A	-15	GLY	-	expression tag	UNP Q9I4X1
A	-14	HIS	-	expression tag	UNP Q9I4X1
A	-13	HIS	-	expression tag	UNP Q9I4X1
A	-12	HIS	-	expression tag	UNP Q9I4X1
A	-11	HIS	-	expression tag	UNP Q9I4X1
A	-10	HIS	-	expression tag	UNP Q9I4X1
A	-9	HIS	-	expression tag	UNP Q9I4X1
A	-8	ALA	-	expression tag	UNP Q9I4X1
A	-7	GLU	-	expression tag	UNP Q9I4X1
A	-6	ASN	-	expression tag	UNP Q9I4X1
A	-5	LEU	-	expression tag	UNP Q9I4X1
A	-4	TYR	-	expression tag	UNP Q9I4X1
A	-3	PHE	-	expression tag	UNP Q9I4X1
A	-2	GLN	-	expression tag	UNP Q9I4X1
A	-1	GLY	-	expression tag	UNP Q9I4X1
A	0	HIS	-	expression tag	UNP Q9I4X1
C	-16	MET	-	initiating methionine	UNP Q9I4X1
C	-15	GLY	-	expression tag	UNP Q9I4X1
C	-14	HIS	-	expression tag	UNP Q9I4X1
C	-13	HIS	-	expression tag	UNP Q9I4X1
C	-12	HIS	-	expression tag	UNP Q9I4X1
C	-11	HIS	-	expression tag	UNP Q9I4X1
C	-10	HIS	-	expression tag	UNP Q9I4X1
C	-9	HIS	-	expression tag	UNP Q9I4X1

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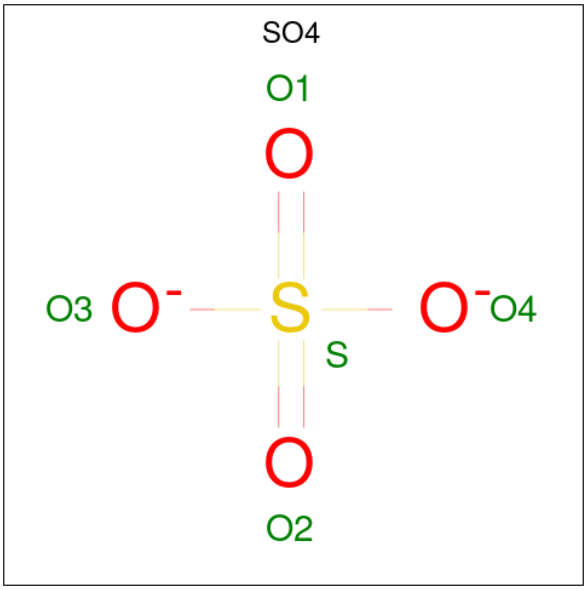
Chain	Residue	Modelled	Actual	Comment	Reference
C	-8	ALA	-	expression tag	UNP Q9I4X1
C	-7	GLU	-	expression tag	UNP Q9I4X1
C	-6	ASN	-	expression tag	UNP Q9I4X1
C	-5	LEU	-	expression tag	UNP Q9I4X1
C	-4	TYR	-	expression tag	UNP Q9I4X1
C	-3	PHE	-	expression tag	UNP Q9I4X1
C	-2	GLN	-	expression tag	UNP Q9I4X1
C	-1	GLY	-	expression tag	UNP Q9I4X1
C	0	HIS	-	expression tag	UNP Q9I4X1

- Molecule 2 is a protein called PqsB.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	279	Total	C	H	N	O	S	0	15	0
			4459	1400	2260	385	405	9			
2	D	283	Total	C	H	N	O	S	0	17	0
			4542	1423	2296	399	415	9			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

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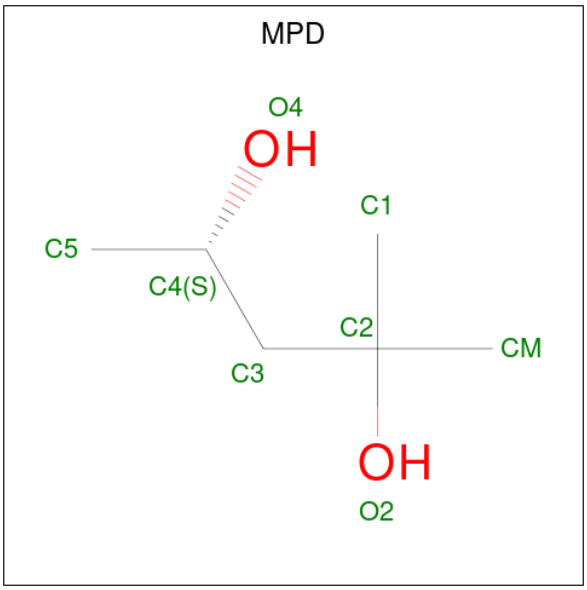
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C H O 22 6 14 2	0	0
4	A	1	Total C H O 22 6 14 2	0	0

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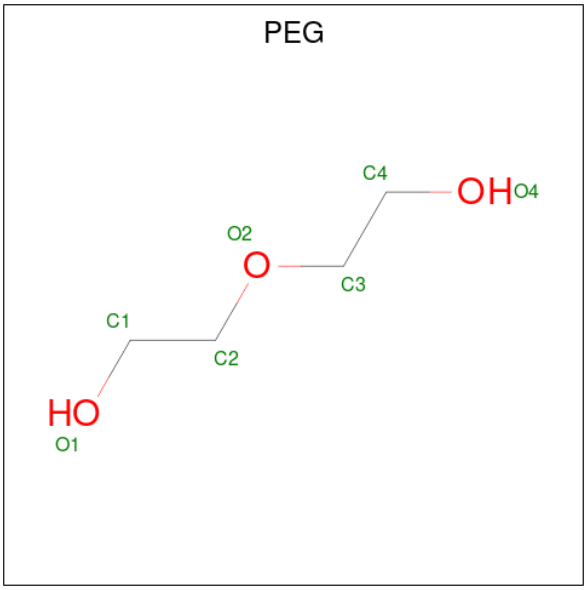
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total C H O 22 6 14 2	0	0
4	C	1	Total C H O 22 6 14 2	0	0
4	C	1	Total C H O 22 6 14 2	0	0

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



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

- Molecule 6 is water.

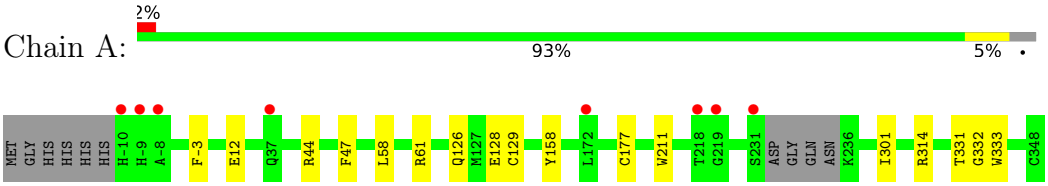
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	361	Total O 361 361	0	0
6	B	271	Total O 271 271	0	0
6	C	282	Total O 282 282	0	0
6	D	298	Total O 298 298	0	1

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.

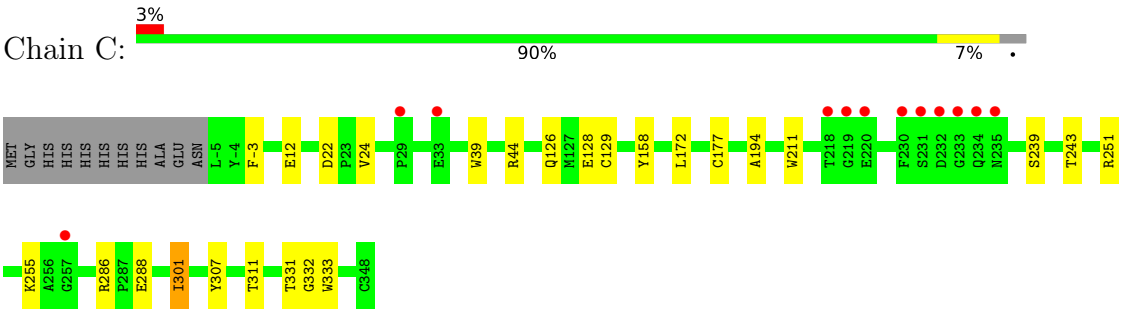
- Molecule 1: PqsC

Chain A:



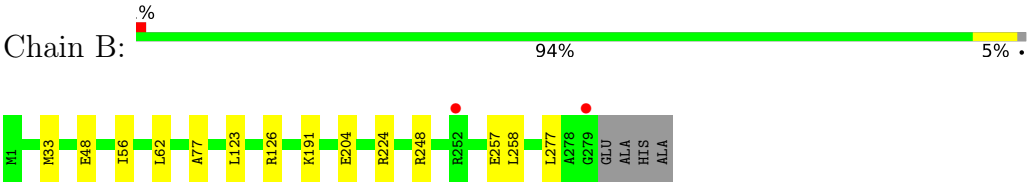
- Molecule 1: PqsC

Chain C:



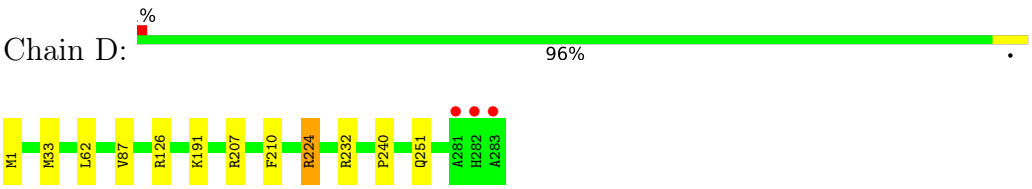
- Molecule 2: PqsB

Chain B:



- Molecule 2: PqsB

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	61.99Å 142.38Å 171.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.34 – 1.84 47.34 – 1.84	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.34-1.84) 99.9 (47.34-1.84)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.25 (at 1.84Å)	Xtriage
Refinement program	PHENIX dev_2875	Depositor
R, R_{free}	0.149 , 0.179 0.150 , 0.179	Depositor DCC
R_{free} test set	6458 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å^2)	19.4	Xtriage
Anisotropy	0.973	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 54.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	21569	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.37% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, PEG, MPD, CSD

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	0.41	0/2901	0.56	0/3936
1	C	0.39	0/2874	0.53	0/3901
2	B	0.40	0/2290	0.56	0/3108
2	D	0.42	0/2340	0.58	0/3173
All	All	0.40	0/10405	0.56	0/14118

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

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	2814	2805	2807	16	0
1	C	2785	2770	2774	18	0
2	B	2199	2260	2259	14	0
2	D	2246	2296	2296	15	0
3	A	20	0	0	1	0
3	B	15	0	0	1	0
3	C	5	0	0	0	0
3	D	15	0	0	0	0
4	A	16	28	28	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	24	42	42	1	0
5	A	7	10	10	0	0
6	A	361	0	0	6	0
6	B	271	0	0	8	1
6	C	282	0	0	2	0
6	D	298	0	0	7	1
All	All	11358	10211	10216	58	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 3.

The worst 5 of 58 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)
2:D:126[A]:ARG:NH2	6:D:1001:HOH:O	1.99	0.93
2:B:126[A]:ARG:NH2	6:B:402:HOH:O	2.03	0.90
1:A:126[B]:GLN:OE1	6:A:1001:HOH:O	1.93	0.86
1:C:286:ARG:NH1	1:C:288:GLU:OE2	2.09	0.85
1:C:126[B]:GLN:NE2	6:C:1001:HOH:O	2.11	0.84

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)
6:B:628:HOH:O	6:D:1214:HOH:O[2_455]	1.97	0.23

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	362/365 (99%)	350 (97%)	11 (3%)	1 (0%)	41 27

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	C	364/365 (100%)	352 (97%)	11 (3%)	1 (0%)	41 27
2	B	293/283 (104%)	289 (99%)	4 (1%)	0	100 100
2	D	298/283 (105%)	294 (99%)	4 (1%)	0	100 100
All	All	1317/1296 (102%)	1285 (98%)	30 (2%)	2 (0%)	47 33

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	301	ILE
1	C	301	ILE

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	301/299 (101%)	299 (99%)	2 (1%)	84 78
1	C	296/299 (99%)	294 (99%)	2 (1%)	84 78
2	B	236/223 (106%)	235 (100%)	1 (0%)	91 88
2	D	239/223 (107%)	235 (98%)	4 (2%)	60 47
All	All	1072/1044 (103%)	1063 (99%)	9 (1%)	81 75

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	224[A]	ARG
2	D	224[B]	ARG
1	C	44	ARG
1	C	158	TYR
2	D	1	MET

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)

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	CSD	C	129[B]	-	3,7,8	0.94	0	1,8,10	0.02	0
1	CSD	C	129[A]	-	3,7,8	0.79	0	1,8,10	0.33	0
1	CSD	A	129[A]	-	3,7,8	0.99	0	1,8,10	0.21	0
1	CSD	A	129[B]	-	3,7,8	0.80	0	1,8,10	1.17	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
1	CSD	C	129[B]	-	-	0/2/6/8	-
1	CSD	C	129[A]	-	-	1/2/6/8	-
1	CSD	A	129[A]	-	-	0/2/6/8	-
1	CSD	A	129[B]	-	-	1/2/6/8	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	129[B]	CSD	CA-CB-SG-OD1
1	C	129[A]	CSD	CA-CB-SG-OD1

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	129[B]	CSD	1	0
1	A	129[B]	CSD	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

17 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	$\#\|Z\| > 2$
4	MPD	C	903	-	7,7,7	0.31	0	9,10,10	0.67	0
3	SO4	A	903	-	4,4,4	0.14	0	6,6,6	0.07	0
3	SO4	B	303	-	4,4,4	0.14	0	6,6,6	0.09	0
4	MPD	A	906	-	7,7,7	0.31	0	9,10,10	0.52	0
3	SO4	A	904	-	4,4,4	0.16	0	6,6,6	0.11	0
3	SO4	A	902	-	4,4,4	0.14	0	6,6,6	0.06	0
3	SO4	D	901	-	4,4,4	0.13	0	6,6,6	0.14	0
4	MPD	C	901	-	7,7,7	0.28	0	9,10,10	0.22	0
3	SO4	A	901	-	4,4,4	0.15	0	6,6,6	0.05	0
3	SO4	B	302	-	4,4,4	0.15	0	6,6,6	0.18	0
3	SO4	D	902	-	4,4,4	0.14	0	6,6,6	0.12	0
4	MPD	C	902	-	7,7,7	0.32	0	9,10,10	0.39	0
5	PEG	A	907	-	6,6,6	0.48	0	5,5,5	0.30	0
4	MPD	A	905	-	7,7,7	0.34	0	9,10,10	0.22	0
3	SO4	C	900	-	4,4,4	0.15	0	6,6,6	0.16	0
3	SO4	B	301	-	4,4,4	0.17	0	6,6,6	0.16	0
3	SO4	D	900	-	4,4,4	0.13	0	6,6,6	0.29	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	MPD	C	903	-	-	3/5/5/5	-
4	MPD	A	906	-	-	0/5/5/5	-
5	PEG	A	907	-	-	1/4/4/4	-
4	MPD	C	901	-	-	0/5/5/5	-
4	MPD	C	902	-	-	0/5/5/5	-
4	MPD	A	905	-	-	1/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	903	MPD	C1-C2-C3-C4
4	C	903	MPD	O2-C2-C3-C4
5	A	907	PEG	C1-C2-O2-C3
4	C	903	MPD	CM-C2-C3-C4
4	A	905	MPD	C2-C3-C4-O4

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	903	MPD	1	0
4	A	906	MPD	1	0
3	A	904	SO4	1	0
3	B	301	SO4	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, 95^{th} 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	354/365 (96%)	-0.39	8 (2%) 60 58	15, 24, 54, 123	0
1	C	353/365 (96%)	-0.22	12 (3%) 45 41	15, 32, 63, 114	0
2	B	279/283 (98%)	-0.38	2 (0%) 87 87	15, 23, 50, 82	0
2	D	283/283 (100%)	-0.41	3 (1%) 80 80	15, 22, 46, 81	0
All	All	1269/1296 (97%)	-0.34	25 (1%) 65 64	15, 25, 56, 123	0

The worst 5 of 25 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	-10	HIS	9.1
1	C	231	SER	6.8
1	A	-9	HIS	6.3
1	C	232	ASP	6.2
1	C	235	ASN	4.9

6.2 Non-standard residues in protein, DNA, RNA chains

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, 95^{th} 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	CSD	C	129[A]	8/9	0.97	0.08	18,24,34,47	0
1	CSD	C	129[B]	8/9	0.97	0.08	18,25,38,46	10
1	CSD	A	129[A]	8/9	0.98	0.08	17,21,29,35	10
1	CSD	A	129[B]	8/9	0.98	0.08	17,19,25,33	0

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, 95^{th} 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	SO4	B	302	5/5	0.83	0.18	74,74,81,88	5
5	PEG	A	907	7/7	0.86	0.14	47,64,78,91	0
3	SO4	A	901	5/5	0.87	0.28	84,86,92,96	0
3	SO4	D	902	5/5	0.87	0.22	81,81,82,95	5
3	SO4	A	904	5/5	0.87	0.17	56,57,61,81	5
4	MPD	A	906	8/8	0.88	0.15	55,66,75,82	0
3	SO4	C	900	5/5	0.89	0.24	36,37,53,58	5
3	SO4	A	903	5/5	0.90	0.13	81,82,88,99	5
4	MPD	C	903	8/8	0.91	0.19	74,90,99,102	0
3	SO4	A	902	5/5	0.93	0.13	60,60,70,76	5
3	SO4	D	901	5/5	0.93	0.13	47,52,58,65	5
3	SO4	B	303	5/5	0.94	0.18	53,55,72,76	5
4	MPD	C	902	8/8	0.94	0.13	37,47,56,67	0
3	SO4	B	301	5/5	0.95	0.10	48,49,53,78	0
4	MPD	A	905	8/8	0.96	0.12	34,43,64,64	0
3	SO4	D	900	5/5	0.96	0.11	34,39,49,55	5
4	MPD	C	901	8/8	0.96	0.08	29,40,69,69	0

6.5 Other polymers (i)

There are no such residues in this entry.