



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 14, 2017 – 04:01 am GMT

PDB ID : 1GKF
Title : Crystal structures of penicillin acylase enzyme-substrate complexes: Structural insights into the catalytic mechanism
Authors : McVey, C.E.; Walsh, M.A.; Dodson, G.G.; Wilson, K.S.; Brannigan, J.A.
Deposited on : 2001-08-13
Resolution : 1.41 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

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

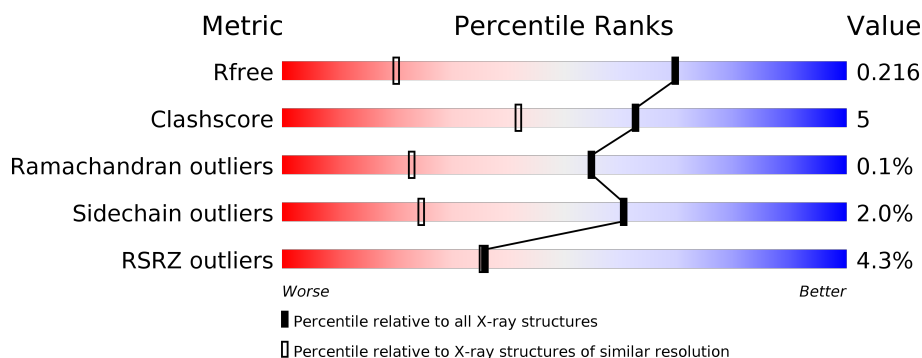
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.41 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}	100719	1836 (1.44-1.40)
Clashscore	112137	1955 (1.44-1.40)
Ramachandran outliers	110173	1908 (1.44-1.40)
Sidechain outliers	110143	1907 (1.44-1.40)
RSRZ outliers	101464	1845 (1.44-1.40)

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. 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	<div> <div>2%</div> <div> <div></div> <div>67%</div> <div>11%</div> <div>••</div> <div>20%</div> </div> </div>
2	B	557	<div> <div>5%</div> <div> <div></div> <div>84%</div> <div>15%</div> <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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EDO	B	1559[A]	-	-	-	X
3	EDO	B	1559[B]	-	-	-	X
3	EDO	B	1560	-	-	-	X
3	EDO	B	1562	-	-	-	X
3	EDO	B	1566	-	-	-	X
3	EDO	B	1569	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7283 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 PENICILLIN G ACYLASE ALPHA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	5	13	1
			1764	1121	300	333	10			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	SME	MET	MODIFIED RESIDUE	UNP P06875

- Molecule 2 is a protein called PENICILLIN G ACYLASE BETA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	557	Total	C	N	O	S	15	18	0
			4562	2896	788	867	11			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	241	ALA	ASN	ENGINEERED MUTATION	UNP P06875

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	1
			7	4	3		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	Ca 1	0	0

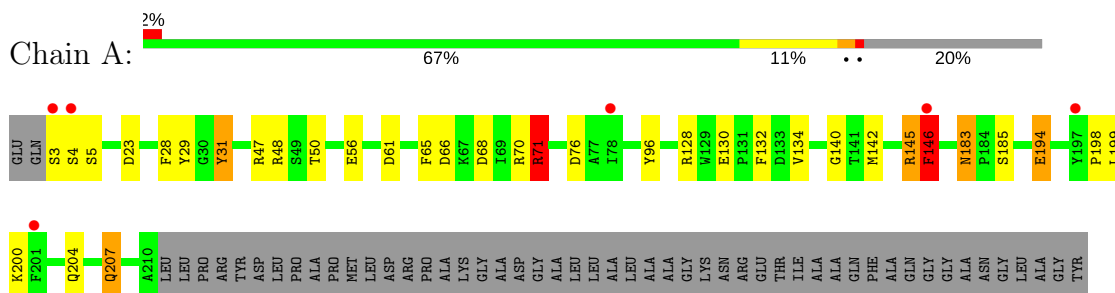
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	241	Total 241	O 241	2	5
5	B	660	Total 660	O 660	0	10

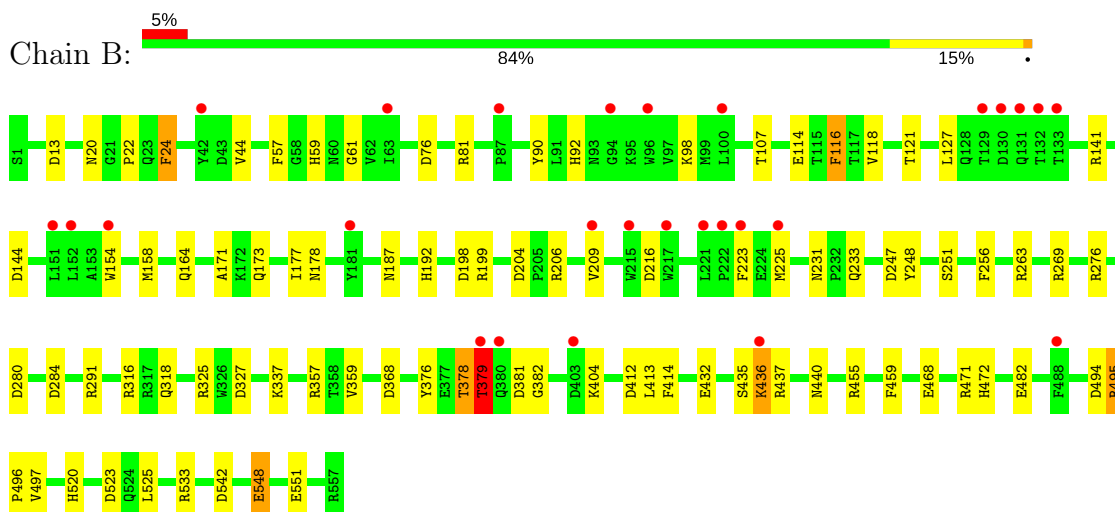
3 Residue-property plots [i](#)

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: PENICILLIN G ACYLASE ALPHA SUBUNIT



• Molecule 2: PENICILLIN G ACYLASE BETA SUBUNIT



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.20Å 131.70Å 63.90Å 90.00° 105.70° 90.00°	Depositor
Resolution (Å)	25.00 – 1.41 24.91 – 1.41	Depositor EDS
% Data completeness (in resolution range)	97.4 (25.00-1.41) 97.2 (24.91-1.41)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.83 (at 1.41Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.148 , 0.168 0.203 , 0.216	Depositor DCC
R_{free} test set	4577 reflections (3.01%)	DCC
Wilson B-factor (Å ²)	10.9	Xtriage
Anisotropy	0.178	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7283	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.39% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, SME, EDO

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	0.96	1/1797 (0.1%)	1.64	33/2435 (1.4%)
2	B	0.98	4/4691 (0.1%)	1.52	55/6396 (0.9%)
All	All	0.97	5/6488 (0.1%)	1.55	88/8831 (1.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
2	B	0	3
All	All	0	6

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	98	LYS	CD-CE	18.66	1.98	1.51
2	B	81	ARG	CD-NE	-11.34	1.27	1.46
1	A	200	LYS	CG-CD	-8.39	1.24	1.52
2	B	432	GLU	CG-CD	-6.50	1.42	1.51
2	B	482	GLU	CD-OE2	5.77	1.31	1.25

The worst 5 of 88 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	455	ARG	NE-CZ-NH2	-18.45	111.08	120.30
2	B	533	ARG	NE-CZ-NH1	17.12	128.86	120.30
2	B	471	ARG	NE-CZ-NH1	13.79	127.19	120.30
1	A	48	ARG	NE-CZ-NH1	13.75	127.18	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	199	ARG	NE-CZ-NH2	-12.82	113.89	120.30

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	146[B]	PHE	Mainchain
1	A	5[B]	SER	Mainchain
1	A	71[B]	ARG	Mainchain
2	B	318[B]	GLN	Mainchain
2	B	376	TYR	Mainchain

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	1764	0	1702	29	0
2	B	4562	0	4364	45	0
3	B	55	0	80	5	0
4	B	1	0	0	0	0
5	A	241	0	0	6	0
5	B	660	0	0	10	2
All	All	7283	0	6146	63	2

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.

The worst 5 of 63 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:B:495[A]:ARG:NH2	5:B:2577[A]:HOH:O	1.70	1.24
2:B:114[A]:GLU:OE1	5:B:2154[A]:HOH:O	1.81	0.98
1:A:207:GLN:HE21	1:A:207:GLN:H	1.18	0.91
2:B:59:HIS:HD2	2:B:61:GLY:H	1.22	0.83
1:A:146[B]:PHE:HB2	5:A:2151:HOH:O	1.79	0.82

All (2) 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:2126:HOH:O	5:B:2508:HOH:O[1_656]	2.04	0.16
5:B:2170:HOH:O	5:B:2378:HOH:O[1_656]	2.17	0.03

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	218/260 (84%)	213 (98%)	5 (2%)	0	100	100
2	B	573/557 (103%)	558 (97%)	14 (2%)	1 (0%)	51	21
All	All	791/817 (97%)	771 (98%)	19 (2%)	1 (0%)	55	23

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	251	SER

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	188/210 (90%)	182 (97%)	6 (3%)	44	10
2	B	477/459 (104%)	469 (98%)	8 (2%)	66	32
All	All	665/669 (99%)	651 (98%)	14 (2%)	60	22

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	20	ASN
2	B	154	TRP
2	B	378[B]	THR
1	A	207	GLN
2	B	378[A]	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 20 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	134	GLN
2	B	187	ASN
2	B	440	ASN
2	B	92	HIS
2	B	93	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	SME	A	16	1	8,8,9	1.24	2 (25%)	6,9,11	1.08	1 (16%)

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	SME	A	16	1	-	0/5/7/9	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	16	SME	OE-S	-2.13	1.44	1.50
1	A	16	SME	CA-C	2.10	1.53	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	16	SME	O-C-CA	-2.27	118.75	125.02

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 1 is monoatomic - leaving 14 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	EDO	B	1558	-	3,3,3	0.58	0	2,2,2	0.21	0
3	EDO	B	1559[A]	-	3,3,3	0.53	0	2,2,2	0.50	0
3	EDO	B	1559[B]	-	3,3,3	0.55	0	2,2,2	0.34	0
3	EDO	B	1560	-	3,3,3	0.64	0	2,2,2	0.18	0
3	EDO	B	1561	-	3,3,3	0.60	0	2,2,2	0.21	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	B	1562	-	3,3,3	0.42	0	2,2,2	1.32	0
3	EDO	B	1563	-	3,3,3	0.57	0	2,2,2	1.36	0
3	EDO	B	1564	-	3,3,3	0.27	0	2,2,2	0.39	0
3	EDO	B	1565	-	3,3,3	0.53	0	2,2,2	0.46	0
3	EDO	B	1566	-	3,3,3	0.60	0	2,2,2	0.48	0
3	EDO	B	1567	-	3,3,3	0.53	0	2,2,2	0.85	0
3	EDO	B	1568	-	3,3,3	0.85	0	2,2,2	2.17	1 (50%)
3	EDO	B	1569	-	3,3,3	1.39	0	2,2,2	0.73	0
3	EDO	B	1570	-	3,3,3	0.64	0	2,2,2	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
3	EDO	B	1558	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1559[A]	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1559[B]	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1560	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1561	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1562	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1563	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1564	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1565	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1566	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1567	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1568	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1569	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1570	-	-	0/1/1/1	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1568	EDO	O1-C1-C2	2.57	130.47	112.08

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1558	EDO	1	0
3	B	1562	EDO	2	0
3	B	1568	EDO	1	0
3	B	1569	EDO	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 [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	207/260 (79%)	0.63	6 (2%)	52	51	7, 11, 21, 44	2 (0%)
2	B	557/557 (100%)	0.51	27 (4%)	31	30	6, 11, 24, 60	5 (0%)
All	All	764/817 (93%)	0.54	33 (4%)	36	35	6, 11, 23, 60	7 (0%)

The worst 5 of 33 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	133	THR	6.5
2	B	132	THR	6.1
1	A	3	SER	4.7
2	B	379[A]	THR	4.5
2	B	131	GLN	4.2

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q < 0.9
1	SME	A	16	9/10	0.95	0.10	-	9,9,11,11	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	EDO	B	1560	4/4	0.93	0.15	11.77	12,15,16,17	0
3	EDO	B	1562	4/4	0.81	0.22	3.52	20,21,21,23	0
3	EDO	B	1559[B]	4/4	0.80	0.18	2.86	20,21,22,26	3
3	EDO	B	1559[A]	4/4	0.80	0.18	2.77	18,19,20,26	3
3	EDO	B	1566	4/4	0.96	0.15	2.76	18,20,22,23	0
3	EDO	B	1569	4/4	0.80	0.18	2.49	15,22,24,25	0
3	EDO	B	1570	4/4	0.97	0.10	0.78	11,16,24,26	0
3	EDO	B	1564	4/4	0.93	0.11	0.67	12,16,16,17	0
3	EDO	B	1563	4/4	0.89	0.12	-0.07	13,17,19,20	0
3	EDO	B	1568	4/4	0.90	0.14	-0.55	14,24,24,27	0
3	EDO	B	1558	4/4	0.96	0.07	-0.63	8,9,10,10	0
4	CA	B	1571	1/1	0.94	0.11	-0.97	7,7,7,7	0
3	EDO	B	1565	4/4	0.95	0.08	-1.10	13,16,16,16	0
3	EDO	B	1567	4/4	0.73	0.26	-	24,25,26,29	0
3	EDO	B	1561	4/4	0.96	0.10	-	17,17,17,18	0

6.5 Other polymers

There are no such residues in this entry.