



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:30 PM UTC

PDB ID : 6YRS / pdb_00006yrs
Title : Structure of a new variant of GNCA ancestral beta-lactamase
Authors : Gavira, J.A.; Risso, V.; Martinez-Rodriguez, S.; Sanchez-Ruiz, J.M.; Modi, T.; Ozkan, S.B.
Deposited on : 2020-04-20
Resolution : 1.70 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
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

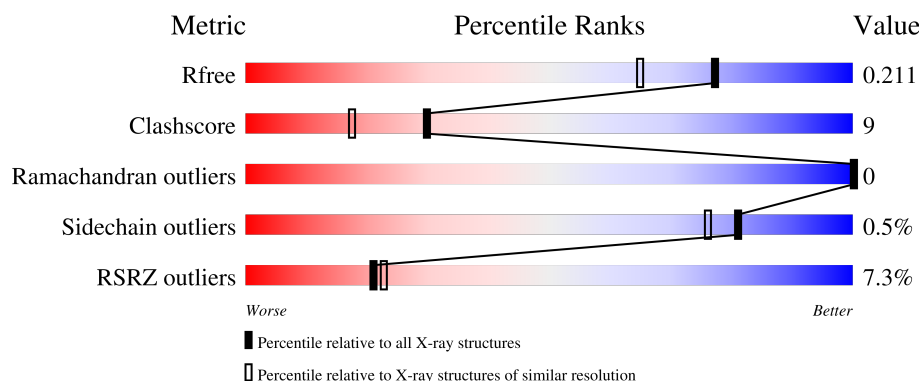
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	269	3% 90% 8% .
1	B	269	11% 71% 17% . 11%

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
6	PEG	B	307	-	-	X	-

2 Entry composition [i](#)

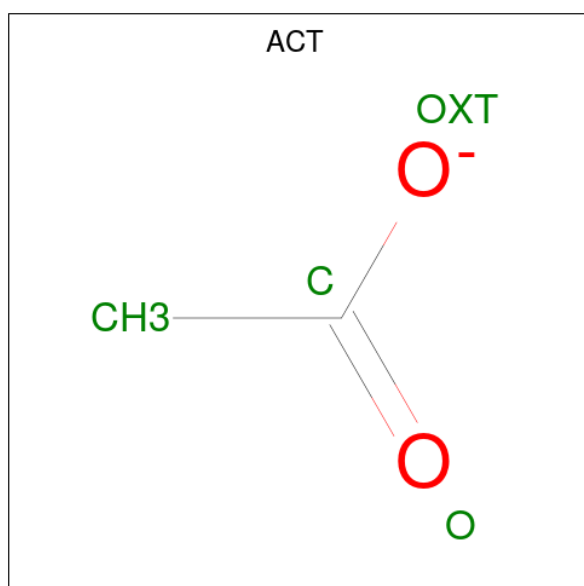
There are 7 unique types of molecules in this entry. The entry contains 8738 atoms, of which 4300 are hydrogens and 0 are deuteriums.

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 ancestral beta-lactamase.

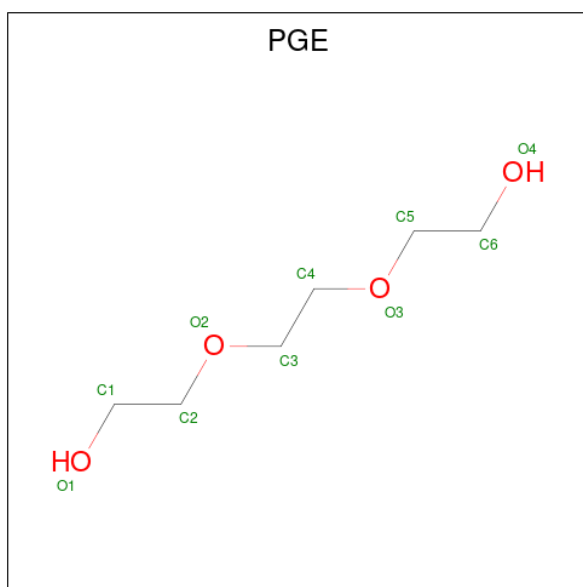
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	264	Total	C	H	N	O	S	0	28	0
			4269	1315	2147	390	411	6			
1	B	240	Total	C	H	N	O	S	0	23	0
			3898	1198	1971	354	368	7			

- Molecule 2 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2^-$).



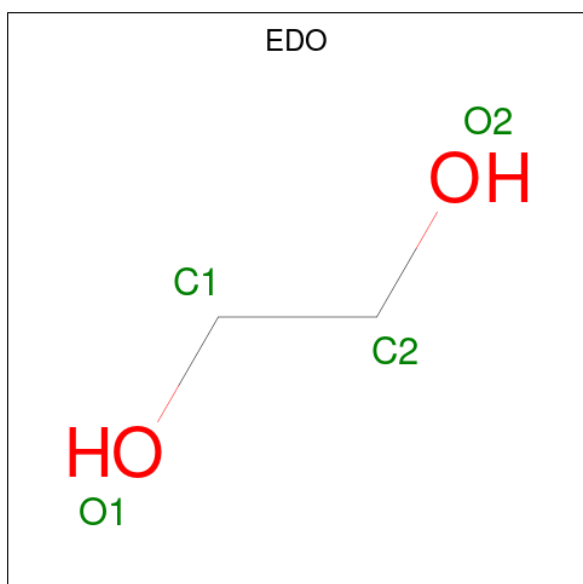
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			7	2	3	2		
2	A	1	Total	C	H	O	0	0
			7	2	3	2		

- Molecule 3 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



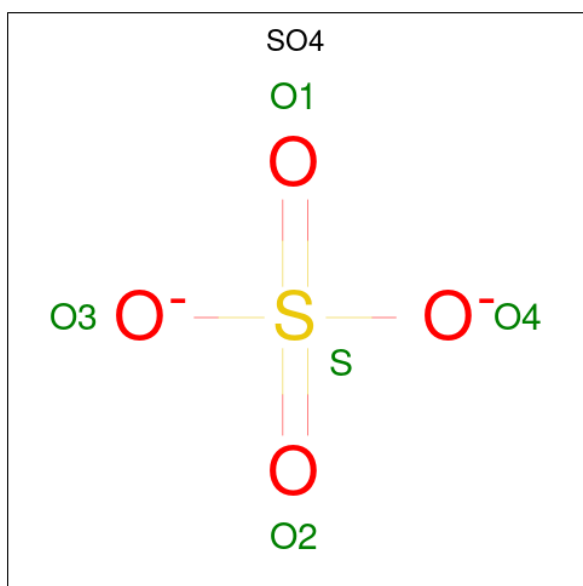
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			24	6	14	4		
3	A	1	Total	C	H	O	0	0
			24	6	14	4		
3	A	1	Total	C	H	O	0	0
			24	6	14	4		
3	A	1	Total	C	H	O	0	1
			48	12	28	8		
3	A	1	Total	C	H	O	0	0
			24	6	14	4		

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



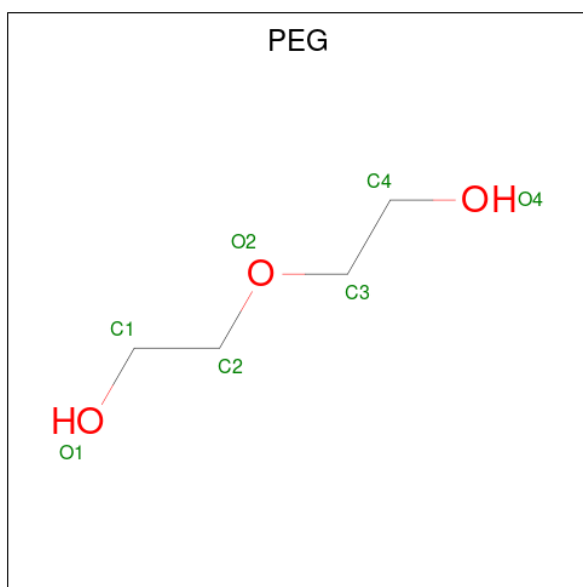
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C H O 10 2 6 2	0	0
4	A	1	Total C H O 10 2 6 2	0	0
4	A	1	Total C H O 10 2 6 2	0	0
4	A	1	Total C H O 10 2 6 2	0	0
4	B	1	Total C H O 10 2 6 2	0	0
4	B	1	Total C H O 10 2 6 2	0	0
4	B	1	Total C H O 10 2 6 2	0	0

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 6 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



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

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	163	Total	O	0	0
			163	163		
7	B	90	Total	O	0	0
			90	90		

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.

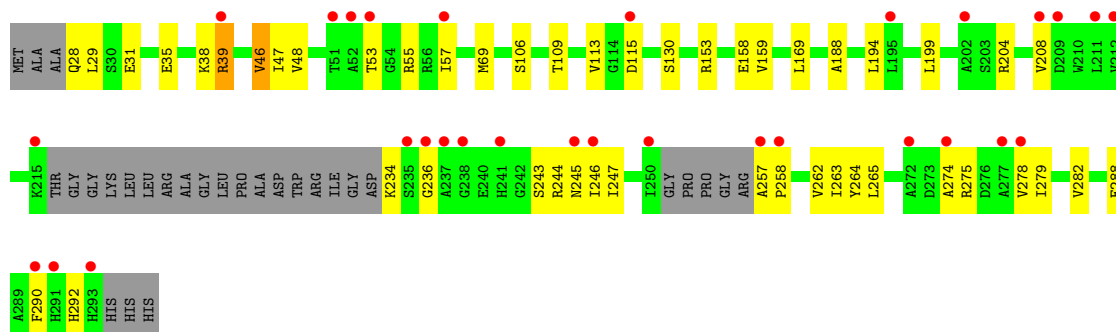
- Molecule 1: ancestral beta-lactamase

Chain A: 



- Molecule 1: ancestral beta-lactamase

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	47.35Å 81.40Å 61.04Å 90.00° 94.58° 90.00°	Depositor
Resolution (Å)	48.73 – 1.70 48.73 – 1.70	Depositor EDS
% Data completeness (in resolution range)	97.7 (48.73-1.70) 97.7 (48.73-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 1.70Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.176 , 0.208 0.179 , 0.211	Depositor DCC
R_{free} test set	2531 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	29.4	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 47.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8738	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.51% 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: SO4, EDO, PEG, PGE, ACT

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	0.56	1/2255 (0.0%)	0.68	0/3057
1	B	0.46	0/2037	0.59	0/2753
All	All	0.51	1/4292 (0.0%)	0.64	0/5810

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	186	MET	CG-SD	5.26	1.94	1.80

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	2122	2147	2040	18	2
1	B	1927	1971	1888	53	1
2	A	8	6	6	0	0
3	A	60	84	84	0	0
4	A	16	24	24	2	0
4	B	12	18	18	1	0
5	A	5	0	0	0	0
6	A	7	10	10	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	28	40	40	6	0
7	A	163	0	0	6	1
7	B	90	0	0	2	0
All	All	4438	4300	4110	72	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 9.

All (72) 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:B:48:VAL:HB	1:B:57:ILE:HG22	1.56	0.85
1:A:267[A]:GLU:OE2	7:A:402:HOH:O	2.02	0.77
1:B:28:GLN:N	1:B:31[A]:GLU:OE2	2.20	0.74
1:B:257:ALA:HB1	1:B:258:PRO:HD3	1.69	0.74
1:A:69:MET:O	1:A:72:PHE:HD1	1.70	0.73
1:A:65[A]:ARG:NH2	7:A:403:HOH:O	2.21	0.73
1:B:48:VAL:HB	1:B:57:ILE:CG2	2.19	0.72
1:B:246:ILE:HD11	1:B:263:ILE:HD12	1.73	0.70
1:A:38:LYS:NZ	7:A:401:HOH:O	1.88	0.69
1:B:188:ALA:HB3	6:B:307:PEG:H41	1.75	0.68
1:B:234[A]:LYS:N	1:B:247:ILE:O	2.27	0.67
1:B:257:ALA:HB1	1:B:258:PRO:CD	2.24	0.67
1:B:69[B]:MET:HG3	1:B:245:ASN:OD1	1.95	0.67
1:B:158:GLU:H	6:B:307:PEG:C1	2.10	0.64
1:B:113:VAL:HG22	7:B:403:HOH:O	1.99	0.62
1:A:256[B]:ARG:NH1	7:A:404:HOH:O	2.32	0.61
1:B:158:GLU:H	6:B:307:PEG:H12	1.66	0.60
1:B:257:ALA:CB	1:B:258:PRO:CD	2.81	0.58
1:B:47:ILE:HB	1:B:262[A]:VAL:HG22	1.86	0.57
1:A:281:GLU:OE1	1:A:284:ARG:NH2	2.33	0.57
1:B:46[B]:VAL:HG21	1:B:282:VAL:CG1	2.35	0.56
1:B:265:LEU:HD22	1:B:278:VAL:HG11	1.87	0.56
1:B:288:GLU:O	1:B:292:HIS:ND1	2.38	0.56
1:B:35:GLU:OE1	1:B:38:LYS:NZ	2.27	0.54
1:B:130:SER:CB	1:B:234[A]:LYS:HZ1	2.21	0.54
1:B:234[B]:LYS:N	1:B:247:ILE:O	2.41	0.53
1:A:202:ALA:HB1	4:A:304:EDO:H22	1.91	0.51
1:B:244:ARG:HG3	1:B:279:ILE:HG13	1.92	0.51
1:A:230[A]:ARG:HH21	1:A:230[A]:ARG:HG3	1.76	0.50
1:B:158:GLU:HB2	6:B:307:PEG:H11	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46[B]:VAL:HG21	1:B:282:VAL:HG11	1.94	0.50
1:B:275:ARG:O	1:B:278:VAL:CG1	2.59	0.50
4:A:309:EDO:H12	1:B:153:ARG:HE	1.77	0.50
1:A:69:MET:O	1:A:72:PHE:CD1	2.60	0.49
1:B:234[A]:LYS:HB3	1:B:247:ILE:HG13	1.94	0.49
1:A:99:LYS:HG2	1:A:113:VAL:HG21	1.95	0.48
1:B:234[A]:LYS:HB3	1:B:247:ILE:CG1	2.43	0.48
1:A:72:PHE:CE1	1:A:166:GLU:OE1	2.66	0.48
1:A:38:LYS:NZ	7:A:408:HOH:O	2.46	0.47
1:A:99:LYS:CE	7:A:429:HOH:O	2.63	0.46
1:A:72:PHE:HE1	1:A:166:GLU:OE1	1.97	0.46
1:B:244:ARG:C	1:B:245:ASN:HD22	2.23	0.46
1:B:35:GLU:OE1	1:B:38:LYS:HD2	2.16	0.45
1:B:69[A]:MET:HE3	1:B:69[A]:MET:HB2	1.83	0.45
1:B:106:SER:HB3	1:B:109:THR:OG1	2.17	0.44
1:B:204:ARG:O	1:B:208:VAL:HG23	2.17	0.44
1:B:39[B]:ARG:NH1	7:B:410:HOH:O	2.52	0.42
1:A:48:VAL:O	1:A:56:ARG:HA	2.19	0.42
1:B:262[B]:VAL:CG1	1:B:264:TYR:CE2	3.02	0.42
4:B:301:EDO:C1	6:B:307:PEG:H32	2.49	0.42
1:A:72:PHE:CZ	1:A:169:LEU:HD21	2.55	0.42
1:B:69[A]:MET:HE1	1:B:243:SER:HB3	2.01	0.42
1:B:194:LEU:HD22	1:B:208:VAL:HG22	2.02	0.42
1:B:29:LEU:HD23	1:B:57:ILE:HD11	2.02	0.42
1:B:46[B]:VAL:CG2	1:B:282:VAL:HG11	2.50	0.41
1:B:47:ILE:HB	1:B:262[A]:VAL:CG2	2.50	0.41
1:B:53:THR:CG2	1:B:55[A]:ARG:HB2	2.50	0.41
1:B:275:ARG:O	1:B:278:VAL:HG12	2.21	0.41
1:B:39[A]:ARG:NE	1:B:39[A]:ARG:HA	2.36	0.41
1:B:55[A]:ARG:NH2	1:B:290:PHE:O	2.52	0.41
1:B:243:SER:HA	1:B:265:LEU:O	2.21	0.41
1:B:159:VAL:HG23	6:B:307:PEG:H22	2.02	0.41
1:B:199:LEU:HB2	1:B:204:ARG:HG3	2.02	0.41
1:B:274:ALA:O	1:B:278:VAL:HG12	2.21	0.41
1:B:169:LEU:C	1:B:169:LEU:HD12	2.46	0.40
1:A:51:THR:HG21	1:A:195:LEU:HD13	2.03	0.40
1:B:247:ILE:HG22	1:B:262[B]:VAL:HG13	2.04	0.40
1:B:275:ARG:C	1:B:278:VAL:HG12	2.47	0.40

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)
1:A:124:GLU:OE2	7:A:401:HOH:O[1_655]	2.07	0.13
1:A:227:ALA:H	1:B:115:ASP:OD2[1_454]	1.47	0.13

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	290/269 (108%)	284 (98%)	6 (2%)	0	100	100
1	B	256/269 (95%)	252 (98%)	4 (2%)	0	100	100
All	All	546/538 (102%)	536 (98%)	10 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	228/206 (111%)	228 (100%)	0	100	100
1	B	210/206 (102%)	206 (98%)	4 (2%)	50	34
All	All	438/412 (106%)	434 (99%)	4 (1%)	81	62

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	39[A]	ARG
1	B	39[B]	ARG

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Mol	Chain	Res	Type
1	B	46[A]	VAL
1	B	46[B]	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	291	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

21 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	$\# Z > 2$
3	PGE	A	307	-	9,9,9	0.30	0	8,8,8	0.33	0
3	PGE	A	306	-	9,9,9	0.33	0	8,8,8	0.25	0
3	PGE	A	303	-	9,9,9	0.33	0	8,8,8	0.32	0
6	PEG	A	313	-	6,6,6	0.50	0	5,5,5	0.35	0
2	ACT	A	302	-	3,3,3	1.15	0	3,3,3	1.62	1 (33%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PGE	A	311	-	9,9,9	0.31	0	8,8,8	0.28	0
6	PEG	B	307	-	6,6,6	0.54	0	5,5,5	0.34	0
3	PGE	A	308[B]	-	9,9,9	0.35	0	8,8,8	0.42	0
4	EDO	B	301	-	3,3,3	0.48	0	2,2,2	0.04	0
4	EDO	B	303	-	3,3,3	0.42	0	2,2,2	0.33	0
4	EDO	A	304	-	3,3,3	0.42	0	2,2,2	0.34	0
4	EDO	A	305	-	3,3,3	0.46	0	2,2,2	0.24	0
6	PEG	B	302	-	6,6,6	0.52	0	5,5,5	0.43	0
6	PEG	B	304	-	6,6,6	0.51	0	5,5,5	0.29	0
2	ACT	A	301	-	3,3,3	1.16	0	3,3,3	1.39	0
3	PGE	A	308[A]	-	9,9,9	0.39	0	8,8,8	0.39	0
4	EDO	A	312	-	3,3,3	0.49	0	2,2,2	0.28	0
5	SO4	A	310	-	4,4,4	0.24	0	6,6,6	0.14	0
6	PEG	B	306	-	6,6,6	0.49	0	5,5,5	0.44	0
4	EDO	B	305	-	3,3,3	0.50	0	2,2,2	0.26	0
4	EDO	A	309	-	3,3,3	0.45	0	2,2,2	0.40	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	EDO	A	309	-	-	1/1/1/1	-
4	EDO	B	301	-	-	1/1/1/1	-
6	PEG	B	306	-	-	2/4/4/4	-
3	PGE	A	311	-	-	4/7/7/7	-
3	PGE	A	307	-	-	7/7/7/7	-
4	EDO	B	305	-	-	1/1/1/1	-
3	PGE	A	306	-	-	3/7/7/7	-
6	PEG	B	302	-	-	2/4/4/4	-
6	PEG	B	307	-	-	2/4/4/4	-
3	PGE	A	308[B]	-	-	4/7/7/7	-
4	EDO	B	303	-	-	1/1/1/1	-
3	PGE	A	303	-	-	2/7/7/7	-
4	EDO	A	304	-	-	1/1/1/1	-
4	EDO	A	305	-	-	1/1/1/1	-
6	PEG	A	313	-	-	3/4/4/4	-
6	PEG	B	304	-	-	1/4/4/4	-
3	PGE	A	308[A]	-	-	2/7/7/7	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	A	312	-	-	1/1/1/1	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	302	ACT	O-C-CH3	-2.00	114.31	122.53

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	308[B]	PGE	O3-C5-C6-O4
6	B	307	PEG	C1-C2-O2-C3
3	A	308[B]	PGE	O2-C3-C4-O3
3	A	308[A]	PGE	O1-C1-C2-O2
3	A	311	PGE	O1-C1-C2-O2
6	B	302	PEG	O2-C3-C4-O4
6	B	306	PEG	O2-C3-C4-O4
6	A	313	PEG	O1-C1-C2-O2
6	B	307	PEG	O2-C3-C4-O4
6	B	304	PEG	O1-C1-C2-O2
4	A	305	EDO	O1-C1-C2-O2
3	A	306	PGE	O1-C1-C2-O2
3	A	303	PGE	O1-C1-C2-O2
4	A	304	EDO	O1-C1-C2-O2
6	B	302	PEG	C4-C3-O2-C2
3	A	307	PGE	O1-C1-C2-O2
3	A	306	PGE	O2-C3-C4-O3
3	A	311	PGE	O3-C5-C6-O4
3	A	308[B]	PGE	C4-C3-O2-C2
3	A	307	PGE	C1-C2-O2-C3
3	A	307	PGE	C3-C4-O3-C5
3	A	307	PGE	O3-C5-C6-O4
3	A	306	PGE	C4-C3-O2-C2
6	A	313	PEG	C1-C2-O2-C3
4	B	305	EDO	O1-C1-C2-O2
6	B	306	PEG	O1-C1-C2-O2
3	A	307	PGE	C4-C3-O2-C2
3	A	308[A]	PGE	O2-C3-C4-O3
4	B	303	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	A	311	PGE	C6-C5-O3-C4
6	A	313	PEG	C4-C3-O2-C2
3	A	307	PGE	C6-C5-O3-C4
4	A	309	EDO	O1-C1-C2-O2
4	A	312	EDO	O1-C1-C2-O2
3	A	303	PGE	C6-C5-O3-C4
4	B	301	EDO	O1-C1-C2-O2
3	A	307	PGE	O2-C3-C4-O3
3	A	311	PGE	O2-C3-C4-O3
3	A	308[B]	PGE	C1-C2-O2-C3

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	307	PEG	6	0
4	B	301	EDO	1	0
4	A	304	EDO	1	0
4	A	309	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

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	264/269 (98%)	-0.01	7 (2%) 56 60	17, 35, 54, 107	15 (5%)
1	B	240/269 (89%)	0.66	30 (12%) 8 8	16, 45, 82, 111	13 (5%)
All	All	504/538 (93%)	0.31	37 (7%) 21 23	16, 39, 74, 111	28 (5%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	236	GLY	4.4
1	B	278	VAL	4.4
1	B	257	ALA	4.1
1	A	294	HIS	4.1
1	B	115	ASP	4.0
1	A	291	HIS	3.9
1	B	291	HIS	3.9
1	A	292	HIS	3.9
1	A	72	PHE	3.8
1	B	293	HIS	3.6
1	B	212	VAL	3.6
1	B	53	THR	3.3
1	B	237	ALA	3.3
1	B	246	ILE	3.1
1	B	208	VAL	3.1
1	B	241	HIS	3.0
1	B	209	ASP	3.0
1	B	238	GLY	3.0
1	A	227	ALA	2.8
1	B	250	ILE	2.8
1	B	274	ALA	2.7
1	B	235[A]	SER	2.6
1	B	290	PHE	2.5
1	A	293	HIS	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	38	LYS	2.4
1	B	57	ILE	2.3
1	B	195	LEU	2.3
1	B	39[A]	ARG	2.3
1	B	215	LYS	2.2
1	B	211	LEU	2.2
1	B	51	THR	2.2
1	B	245	ASN	2.2
1	B	272	ALA	2.1
1	B	258	PRO	2.1
1	B	277	ALA	2.1
1	B	52	ALA	2.0
1	B	202	ALA	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(Å ²)	Q<0.9
3	PGE	A	308[A]	10/10	0.66	0.23	37,47,57,57	24
3	PGE	A	308[B]	10/10	0.66	0.23	44,54,56,57	24
6	PEG	A	313	7/7	0.76	0.18	81,97,100,100	0
6	PEG	B	306	7/7	0.77	0.25	70,84,91,91	0
4	EDO	A	312	4/4	0.81	0.19	83,100,102,102	0
4	EDO	B	305	4/4	0.81	0.18	70,84,87,87	0
2	ACT	A	302	4/4	0.81	0.14	58,58,70,70	0
4	EDO	A	309	4/4	0.81	0.18	77,92,93,93	0
4	EDO	B	301	4/4	0.83	0.12	56,67,68,70	0
6	PEG	B	307	7/7	0.83	0.15	75,90,95,95	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	PGE	A	307	10/10	0.84	0.16	63,76,80,80	0
6	PEG	B	304	7/7	0.85	0.15	77,93,94,94	0
4	EDO	A	305	4/4	0.85	0.12	62,75,77,77	0
6	PEG	B	302	7/7	0.85	0.14	65,78,84,84	0
5	SO4	A	310	5/5	0.86	0.16	95,95,96,96	0
3	PGE	A	311	10/10	0.87	0.17	69,83,89,89	0
3	PGE	A	306	10/10	0.87	0.13	64,77,80,80	0
3	PGE	A	303	10/10	0.88	0.12	59,71,78,79	0
4	EDO	A	304	4/4	0.89	0.13	47,57,59,59	0
4	EDO	B	303	4/4	0.92	0.11	54,65,65,66	0
2	ACT	A	301	4/4	0.94	0.08	48,49,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.