



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:32 PM GMT

PDB ID : 3ZST
Title : GLGE ISOFORM 1 FROM STREPTOMYCES COELICOLOR WITH
ALPHA-CYCLODEXTRIN BOUND
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Deposited on : 2011-06-30
Resolution : 2.30 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
<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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

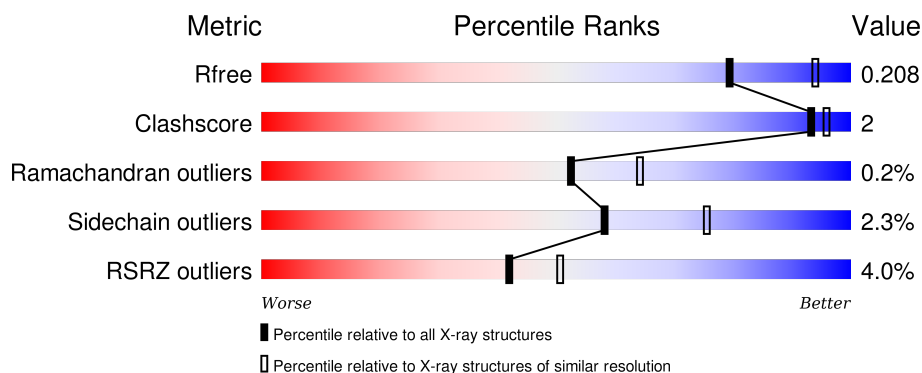
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	695	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 89%, grey 7%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 89% 7% </div> </div>
1	B	695	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 6%, orange 1%, yellow 1%, green 87%, grey 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 6% 87% 5% • 7% </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
2	ACX	A	700	-	-	-	X
2	ACX	B	700	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11175 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 PUTATIVE GLUCANOHYDROLASE PEP1A GLGE ISO-FORM 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	649	Total	C	N	O	S	0	5	0
			5157	3260	937	950	10			
1	B	649	Total	C	N	O	S	0	5	0
			5151	3257	936	948	10			

There are 40 discrepancies between the modelled and reference sequences:

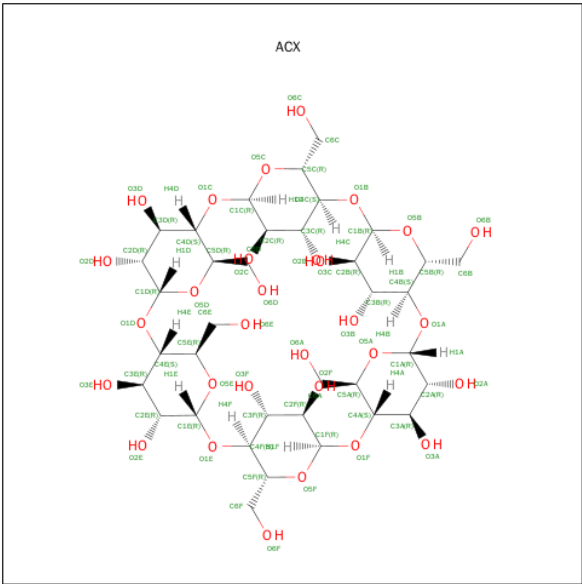
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP Q9L1K2
A	-18	GLY	-	EXPRESSION TAG	UNP Q9L1K2
A	-17	SER	-	EXPRESSION TAG	UNP Q9L1K2
A	-16	SER	-	EXPRESSION TAG	UNP Q9L1K2
A	-15	HIS	-	EXPRESSION TAG	UNP Q9L1K2
A	-14	HIS	-	EXPRESSION TAG	UNP Q9L1K2
A	-13	HIS	-	EXPRESSION TAG	UNP Q9L1K2
A	-12	HIS	-	EXPRESSION TAG	UNP Q9L1K2
A	-11	HIS	-	EXPRESSION TAG	UNP Q9L1K2
A	-10	HIS	-	EXPRESSION TAG	UNP Q9L1K2
A	-9	SER	-	EXPRESSION TAG	UNP Q9L1K2
A	-8	SER	-	EXPRESSION TAG	UNP Q9L1K2
A	-7	GLY	-	EXPRESSION TAG	UNP Q9L1K2
A	-6	LEU	-	EXPRESSION TAG	UNP Q9L1K2
A	-5	VAL	-	EXPRESSION TAG	UNP Q9L1K2
A	-4	PRO	-	EXPRESSION TAG	UNP Q9L1K2
A	-3	ARG	-	EXPRESSION TAG	UNP Q9L1K2
A	-2	GLY	-	EXPRESSION TAG	UNP Q9L1K2
A	-1	SER	-	EXPRESSION TAG	UNP Q9L1K2
A	0	HIS	-	EXPRESSION TAG	UNP Q9L1K2
B	-19	MET	-	EXPRESSION TAG	UNP Q9L1K2
B	-18	GLY	-	EXPRESSION TAG	UNP Q9L1K2
B	-17	SER	-	EXPRESSION TAG	UNP Q9L1K2
B	-16	SER	-	EXPRESSION TAG	UNP Q9L1K2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-15	HIS	-	EXPRESSION TAG	UNP Q9L1K2
B	-14	HIS	-	EXPRESSION TAG	UNP Q9L1K2
B	-13	HIS	-	EXPRESSION TAG	UNP Q9L1K2
B	-12	HIS	-	EXPRESSION TAG	UNP Q9L1K2
B	-11	HIS	-	EXPRESSION TAG	UNP Q9L1K2
B	-10	HIS	-	EXPRESSION TAG	UNP Q9L1K2
B	-9	SER	-	EXPRESSION TAG	UNP Q9L1K2
B	-8	SER	-	EXPRESSION TAG	UNP Q9L1K2
B	-7	GLY	-	EXPRESSION TAG	UNP Q9L1K2
B	-6	LEU	-	EXPRESSION TAG	UNP Q9L1K2
B	-5	VAL	-	EXPRESSION TAG	UNP Q9L1K2
B	-4	PRO	-	EXPRESSION TAG	UNP Q9L1K2
B	-3	ARG	-	EXPRESSION TAG	UNP Q9L1K2
B	-2	GLY	-	EXPRESSION TAG	UNP Q9L1K2
B	-1	SER	-	EXPRESSION TAG	UNP Q9L1K2
B	0	HIS	-	EXPRESSION TAG	UNP Q9L1K2

- Molecule 2 is SUGAR (ALPHA-CYCLODEXTRIN (CYCLOHEXA-AMYLOSE)) (three-letter code: ACX) (formula: C₃₆H₆₀O₃₀).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			66	36	30		
2	B	1	Total	C	O	0	0
			66	36	30		

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



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

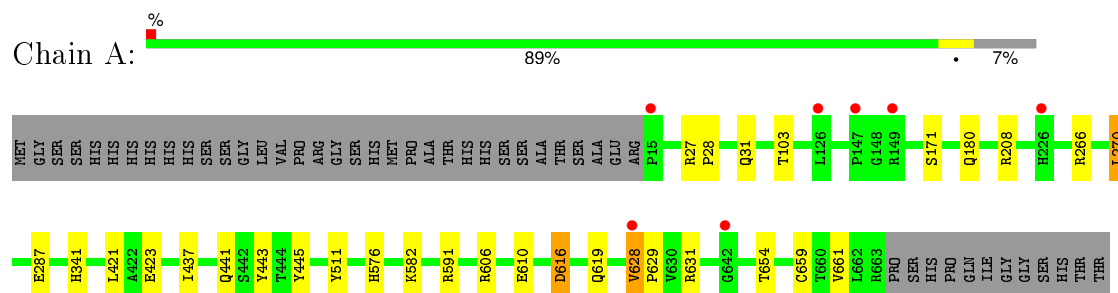
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	488	Total	O	0	0
			488	488		
4	B	243	Total	O	0	0
			243	243		

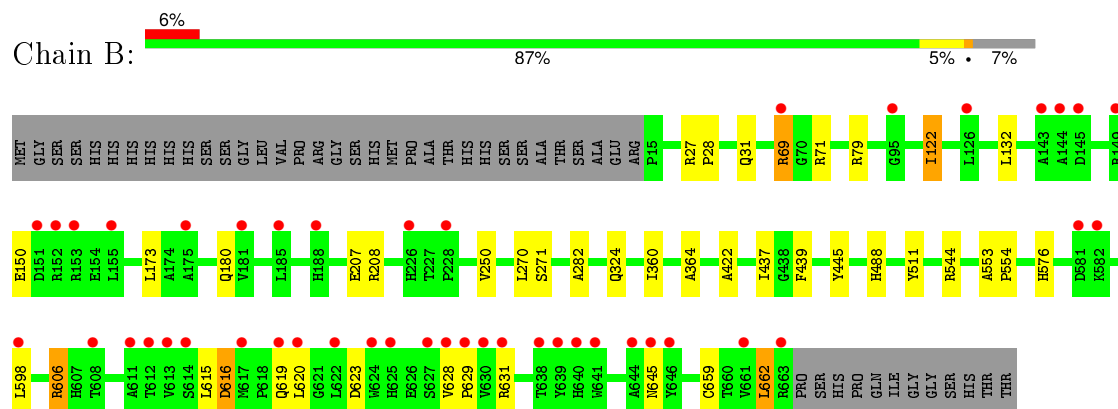
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: PUTATIVE GLUCANOHYDROLASE PEP1A GLGE ISOFORM 1



- Molecule 1: PUTATIVE GLUCANOHYDROLASE PEP1A GLGE ISOFORM 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	113.24Å 113.24Å 314.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	106.54 – 2.30 56.10 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.4 (106.54-2.30) 93.4 (56.10-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.16 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.6.0101	Depositor
R, R_{free}	0.173 , 0.201 0.180 , 0.208	Depositor DCC
R_{free} test set	4305 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	27.2	Xtriage
Anisotropy	0.445	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 31.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 85748 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11175	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.92% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: ACX, 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	0.80	1/5324 (0.0%)	0.78	4/7280 (0.1%)
1	B	0.71	0/5319	0.73	1/7275 (0.0%)
All	All	0.75	1/10643 (0.0%)	0.76	5/14555 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	287	GLU	CD-OE1	5.36	1.31	1.25

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	208	ARG	NE-CZ-NH1	6.63	123.61	120.30
1	B	208	ARG	NE-CZ-NH1	5.74	123.17	120.30
1	A	606	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	A	266[A]	ARG	NE-CZ-NH2	-5.10	117.75	120.30
1	A	266[B]	ARG	NE-CZ-NH2	-5.10	117.75	120.30

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	5157	0	5002	14	0
1	B	5151	0	4990	25	0
2	A	66	0	60	1	0
2	B	66	0	60	1	0
3	A	4	0	6	0	0
4	A	488	0	0	3	0
4	B	243	0	0	1	0
All	All	11175	0	10118	39	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 (39) 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:615:LEU:HD23	1:B:620:LEU:HD11	1.52	0.90
1:B:628:VAL:HG13	1:B:629:PRO:HD2	1.63	0.78
1:B:282:ALA:HA	1:B:360:ILE:HD13	1.69	0.73
1:B:122:ILE:CD1	1:B:132:LEU:HD21	2.19	0.72
1:A:628:VAL:HG13	1:A:629:PRO:HD2	1.77	0.66
1:B:122:ILE:HD11	1:B:132:LEU:HD21	1.78	0.65
1:B:122:ILE:HD11	1:B:132:LEU:CD2	2.35	0.56
1:B:364:ALA:HB1	4:B:2175:HOH:O	2.07	0.54
1:A:103:THR:HG21	4:A:2119:HOH:O	2.07	0.54
1:B:615:LEU:HD23	1:B:620:LEU:CD1	2.33	0.52
1:B:615:LEU:HD13	1:B:645:ASN:ND2	2.26	0.51
1:A:576:HIS:CG	1:A:619:GLN:HG2	2.48	0.49
1:B:576:HIS:CG	1:B:619:GLN:HG2	2.47	0.49
1:B:282:ALA:HA	1:B:360:ILE:CD1	2.40	0.48
1:B:437:ILE:HD12	1:B:437:ILE:C	2.34	0.48
1:A:27:ARG:HB3	1:A:28:PRO:HA	1.97	0.47
1:B:324:GLN:NE2	1:B:360:ILE:HD12	2.31	0.46
1:B:445:TYR:CE1	2:B:700:ACX:H1A	2.51	0.46
1:B:488:HIS:HB3	1:B:606[B]:ARG:HH21	1.81	0.45
1:A:591:ARG:HD3	4:A:2459:HOH:O	2.15	0.45
1:B:422:ALA:HB2	1:B:439:PHE:CD1	2.52	0.45
1:B:553:ALA:HB3	1:B:554:PRO:HD3	1.98	0.44
1:A:445:TYR:CE1	2:A:700:ACX:H1E	2.53	0.44
1:A:437:ILE:HD12	1:A:437:ILE:C	2.38	0.44
1:B:27:ARG:HB3	1:B:28:PRO:HA	1.99	0.44
1:B:662:LEU:H	1:B:662:LEU:HD12	1.83	0.43
1:B:250:VAL:HG12	1:B:250:VAL:O	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:423:GLU:HA	1:A:443:TYR:CD1	2.54	0.42
1:A:270:LEU:HD12	1:A:270:LEU:HA	1.80	0.42
1:A:341:HIS:HE1	4:A:2301:HOH:O	2.02	0.42
1:A:582:LYS:HG3	1:A:610:GLU:O	2.20	0.41
1:B:615:LEU:HD12	1:B:615:LEU:N	2.36	0.41
1:B:628:VAL:HG13	1:B:629:PRO:CD	2.43	0.41
1:A:421:LEU:HD13	1:A:441:GLN:HB3	2.01	0.41
1:B:122:ILE:HD12	1:B:173:LEU:HB2	2.03	0.41
1:A:628:VAL:CG1	1:A:661:VAL:HG13	2.51	0.41
1:B:631:ARG:O	1:B:659:CYS:HA	2.22	0.40
1:A:631:ARG:O	1:A:659:CYS:HA	2.21	0.40
1:B:69:ARG:HB2	1:B:69:ARG:HE	1.74	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	652/695 (94%)	640 (98%)	11 (2%)	1 (0%)	52	64
1	B	652/695 (94%)	641 (98%)	10 (2%)	1 (0%)	52	64
All	All	1304/1390 (94%)	1281 (98%)	21 (2%)	2 (0%)	52	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	616	ASP
1	A	616	ASP

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	539/575 (94%)	531 (98%)	8 (2%)	72	85
1	B	538/575 (94%)	520 (97%)	18 (3%)	45	61
All	All	1077/1150 (94%)	1051 (98%)	26 (2%)	58	74

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	171	SER
1	A	180	GLN
1	A	270	LEU
1	A	511	TYR
1	A	616	ASP
1	A	628	VAL
1	A	654	THR
1	B	31	GLN
1	B	69	ARG
1	B	71	ARG
1	B	79	ARG
1	B	122	ILE
1	B	150	GLU
1	B	180	GLN
1	B	207	GLU
1	B	270	LEU
1	B	271	SER
1	B	511	TYR
1	B	544	ARG
1	B	598	LEU
1	B	606[A]	ARG
1	B	606[B]	ARG
1	B	616	ASP
1	B	623	ASP
1	B	662	LEU

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

Mol	Chain	Res	Type
1	A	262	HIS
1	A	397	HIS
1	A	436	GLN
1	B	262	HIS
1	B	397	HIS
1	B	436	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

3 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	$\# Z > 2$
2	ACX	A	700	-	72,72,72	1.23	6 (8%)	108,108,108	1.57	25 (23%)
3	EDO	A	750	-	3,3,3	0.79	0	2,2,2	0.11	0
2	ACX	B	700	-	72,72,72	1.34	8 (11%)	108,108,108	1.41	15 (13%)

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
2	ACX	A	700	-	-	0/36/156/156	0/0/7/7
3	EDO	A	750	-	-	0/1/1/1	0/0/0/0
2	ACX	B	700	-	-	0/36/156/156	0/0/7/7

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	700	ACX	O1B-C1B	2.04	1.47	1.41
2	A	700	ACX	C4A-C5A	2.05	1.58	1.52
2	B	700	ACX	O5B-C1B	2.08	1.47	1.41
2	A	700	ACX	O1E-C4F	2.10	1.49	1.43
2	B	700	ACX	O5C-C1C	2.11	1.47	1.41
2	A	700	ACX	O2E-C2E	2.32	1.48	1.43
2	B	700	ACX	O5E-C5E	2.42	1.50	1.44
2	A	700	ACX	O1E-C1E	2.51	1.48	1.41
2	B	700	ACX	C4C-C5C	2.56	1.60	1.52
2	A	700	ACX	O5A-C1A	2.58	1.48	1.41
2	B	700	ACX	O1E-C1E	2.65	1.48	1.41
2	B	700	ACX	O5E-C1E	3.01	1.49	1.41
2	A	700	ACX	O5E-C1E	3.17	1.50	1.41
2	B	700	ACX	O5A-C1A	3.64	1.51	1.41

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	700	ACX	O1A-C4B-C3B	-4.27	96.14	107.17
2	B	700	ACX	O1A-C1A-C2A	-4.02	98.32	108.10
2	A	700	ACX	C6F-C5F-C4F	-3.43	103.26	113.25
2	B	700	ACX	C1D-O5D-C5D	-3.07	107.78	113.75
2	A	700	ACX	O5F-C5F-C4F	-2.72	104.00	109.75
2	A	700	ACX	O5F-C1F-C2F	-2.68	104.77	110.28
2	A	700	ACX	O1A-C1A-O5A	-2.54	104.25	110.68
2	A	700	ACX	C1D-O1D-C4E	-2.48	111.53	118.01
2	A	700	ACX	C1E-C2E-C3E	-2.44	105.17	109.97
2	B	700	ACX	O5B-C5B-C4B	-2.40	104.67	109.75
2	A	700	ACX	C6C-C5C-C4C	-2.37	106.35	113.25
2	A	700	ACX	O1C-C1C-O5C	-2.11	105.35	110.68
2	A	700	ACX	O1E-C1E-C2E	-2.07	103.07	108.10
2	B	700	ACX	O5F-C5F-C4F	2.01	114.00	109.75
2	B	700	ACX	C2F-C3F-C4F	2.09	114.18	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	700	ACX	O1C-C1C-C2C	2.09	113.20	108.10
2	A	700	ACX	C3B-C4B-C5B	2.20	115.81	110.84
2	B	700	ACX	O5A-C5A-C6A	2.20	111.93	106.36
2	A	700	ACX	O5B-C5B-C4B	2.21	114.41	109.75
2	A	700	ACX	C3E-C4E-C5E	2.21	115.83	110.84
2	B	700	ACX	O5B-C5B-C6B	2.25	112.04	106.36
2	A	700	ACX	O1B-C4C-C3C	2.26	113.00	107.17
2	B	700	ACX	O5A-C1A-C2A	2.27	114.93	110.28
2	A	700	ACX	C1C-O5C-C5C	2.39	118.39	113.75
2	B	700	ACX	C1E-O5E-C5E	2.50	118.59	113.75
2	A	700	ACX	O5A-C1A-C2A	2.50	115.40	110.28
2	A	700	ACX	C3D-C4D-C5D	2.54	116.58	110.84
2	B	700	ACX	O1A-C1A-O5A	2.55	117.14	110.68
2	B	700	ACX	O1E-C1E-O5E	2.60	117.27	110.68
2	A	700	ACX	O5D-C1D-C2D	2.64	115.69	110.28
2	A	700	ACX	O5A-C5A-C6A	2.74	113.28	106.36
2	B	700	ACX	C3A-C4A-C5A	2.83	117.23	110.84
2	A	700	ACX	O1C-C1C-C2C	2.85	115.04	108.10
2	A	700	ACX	C2E-C3E-C4E	2.91	116.00	109.60
2	A	700	ACX	C3A-C4A-C5A	2.94	117.49	110.84
2	A	700	ACX	O5E-C1E-C2E	3.08	116.59	110.28
2	A	700	ACX	O5D-C5D-C4D	3.15	116.39	109.75
2	B	700	ACX	O5D-C5D-C6D	3.43	115.03	106.36
2	B	700	ACX	O5E-C5E-C4E	3.99	118.18	109.75
2	A	700	ACX	O5C-C5C-C4C	4.55	119.35	109.75

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	700	ACX	1	0
2	B	700	ACX	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	649/695 (93%)	-0.16	7 (1%) 82 86	26, 36, 64, 103	0
1	B	649/695 (93%)	0.23	45 (6%) 20 27	36, 53, 86, 121	0
All	All	1298/1390 (93%)	0.04	52 (4%) 42 51	26, 46, 78, 121	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	144	ALA	6.4
1	B	641	TRP	6.3
1	B	640	HIS	4.7
1	B	625	HIS	4.6
1	B	145	ASP	4.4
1	B	644	ALA	4.4
1	B	628	VAL	4.2
1	B	622	LEU	4.2
1	B	639	TYR	4.1
1	B	661	VAL	4.1
1	B	627	SER	3.9
1	B	181	VAL	3.7
1	B	598	LEU	3.5
1	B	613	VAL	3.5
1	B	630	VAL	3.1
1	B	614	SER	3.1
1	B	69	ARG	3.0
1	A	642	GLY	3.0
1	B	638	THR	2.9
1	B	612	THR	2.8
1	B	151	ASP	2.7
1	B	629	PRO	2.7
1	B	126	LEU	2.7
1	B	620	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	153	ARG	2.7
1	B	149	ARG	2.6
1	B	617	MET	2.6
1	B	226	HIS	2.6
1	B	624	TRP	2.6
1	B	581	ASP	2.5
1	B	619	GLN	2.5
1	A	15	PRO	2.5
1	B	646	TYR	2.5
1	B	611	ALA	2.4
1	B	663	ARG	2.4
1	B	175	ALA	2.4
1	B	155	LEU	2.3
1	A	149	ARG	2.3
1	A	226	HIS	2.3
1	B	228	PRO	2.2
1	B	608	THR	2.2
1	B	143	ALA	2.2
1	A	147	PRO	2.1
1	B	188	HIS	2.1
1	A	126	LEU	2.1
1	A	628	VAL	2.1
1	B	645	ASN	2.1
1	B	185	LEU	2.0
1	B	582	LYS	2.0
1	B	95	GLY	2.0
1	B	631	ARG	2.0
1	B	152	ARG	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 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. 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
2	ACX	B	700	66/66	0.78	0.32	13.36	60,81,108,118	0
2	ACX	A	700	66/66	0.86	0.19	2.94	46,70,100,102	0
3	EDO	A	750	4/4	0.91	0.15	1.55	42,46,52,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.