



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:33 AM UTC

PDB ID : 4BA0 / pdb_00004ba0
Title : Crystal Structure of Agd31B, alpha-transglucosylase, complexed with 5F-alpha-GlcF
Authors : Larsbrink, J.; Izumi, A.; Hemsworth, G.R.; Davies, G.J.; Brumer, H.
Deposited on : 2012-09-09
Resolution : 1.85 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

<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

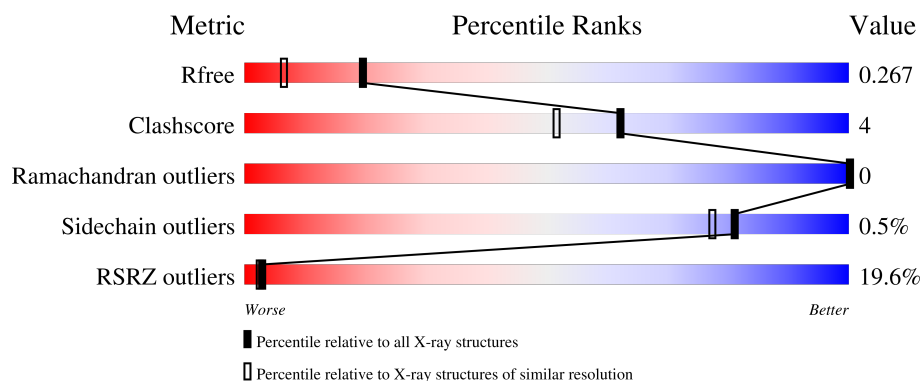
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	817	

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
7	PEG	A	1828	-	-	X	-

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 6830 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.

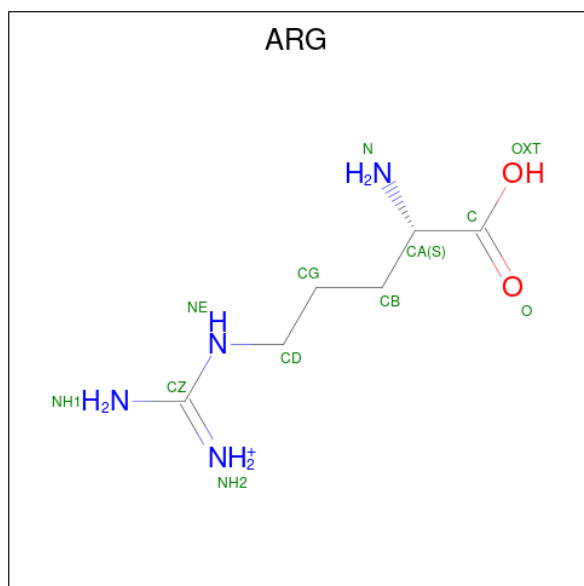
- Molecule 1 is a protein called ALPHA-GLUCOSIDASE, PUTATIVE, ADG31B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	781	6193	3968	1049	1153	23	0	5	0

There is a discrepancy between the modelled and reference sequences:

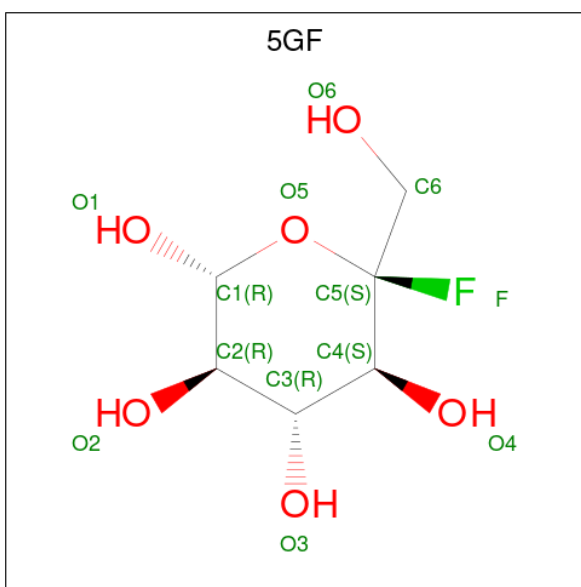
Chain	Residue	Modelled	Actual	Comment	Reference
A	817	LYS	-	expression tag	UNP B3PEE6

- Molecule 2 is ARGinine (CCD ID: ARG) (formula: $C_6H_{15}N_4O_2$).



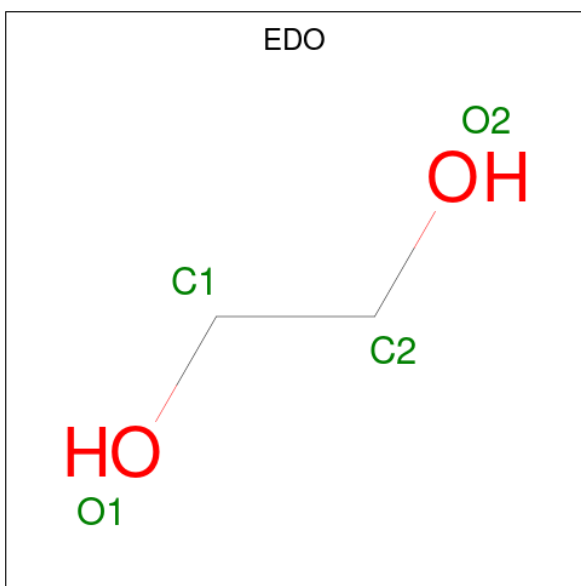
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	12	6	4	2	0	0

- Molecule 3 is 5-fluoro-beta-D-glucopyranose (CCD ID: 5GF) (formula: $C_6H_{11}FO_6$).



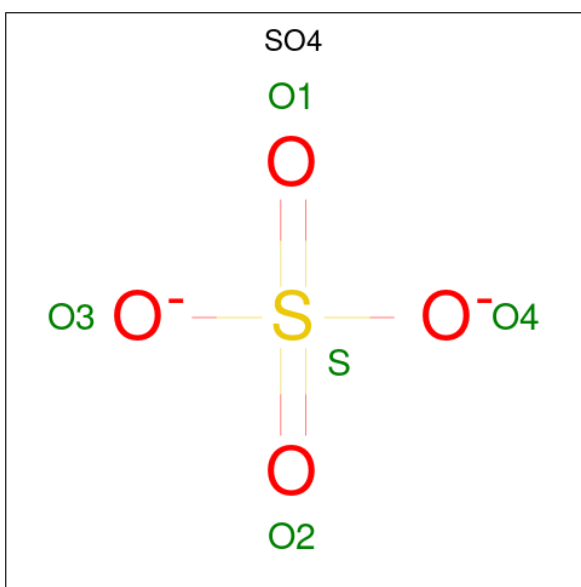
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	F	O	0	0
			12	6	1	5		

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



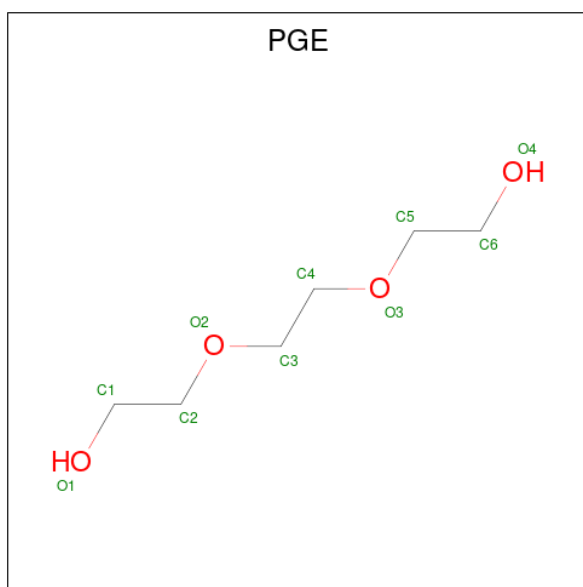
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



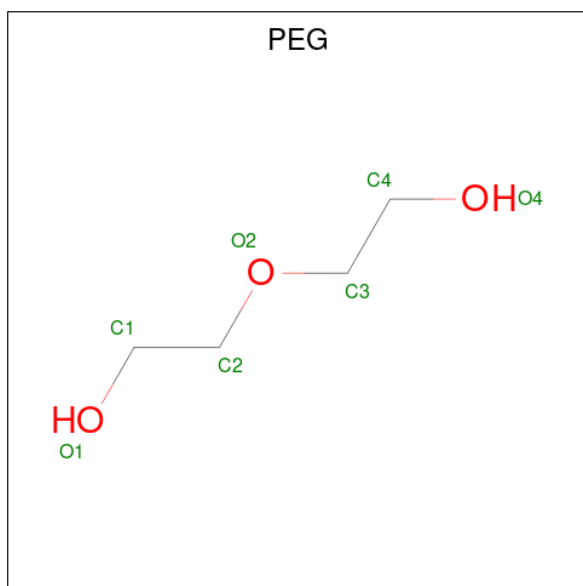
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



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

- Molecule 7 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

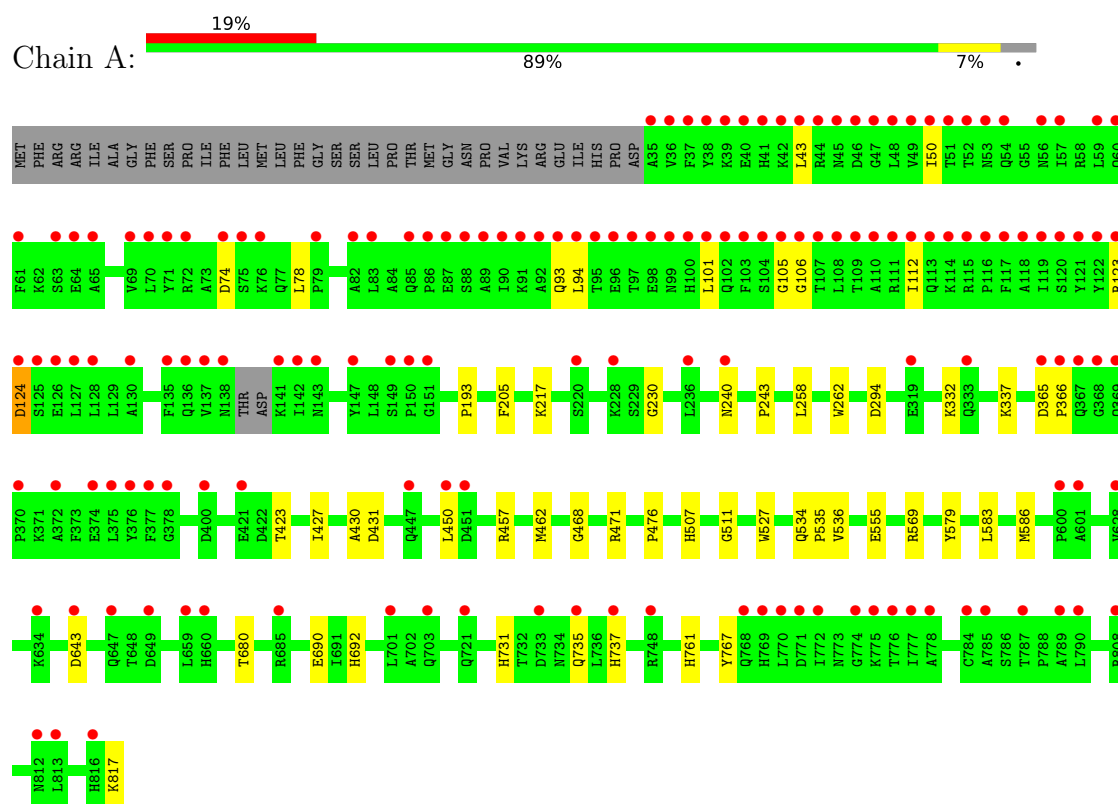
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	557	Total 557	O 557	0	0

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: ALPHA-GLUCOSIDASE, PUTATIVE, ADG31B



4 Data and refinement statistics

Property	Value	Source
Space group	P 6 2 2	Depositor
Cell constants a, b, c, α , β , γ	197.23Å 197.23Å 103.06Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.48 – 1.85 44.48 – 1.85	Depositor EDS
% Data completeness (in resolution range)	100.0 (44.48-1.85) 100.0 (44.48-1.85)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.97 (at 1.86Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.167 , 0.193 0.262 , 0.267	Depositor DCC
R_{free} test set	5007 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	0.849	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.2	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	0.93	EDS
Total number of atoms	6830	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.92% 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, 0.333 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: 5GF, EDO, PEG, PGE, SO4

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	0/6368	0.72	1/8645 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	124	ASP	CB-CA-C	-5.40	109.89	117.23

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	6193	0	5966	50	0
2	A	12	0	12	0	0
3	A	12	0	9	0	0
4	A	4	0	6	1	0
5	A	35	0	0	0	0
6	A	10	0	14	3	0
7	A	7	0	10	7	0
8	A	557	0	0	5	0
All	All	6830	0	6017	50	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 4.

All (50) 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:A:692:HIS:CD2	1:A:761:HIS:HE1	1.78	1.01
1:A:569:ARG:NH2	6:A:1827:PGE:H12	1.78	0.98
1:A:692:HIS:HD2	1:A:761:HIS:HE1	1.12	0.96
1:A:692:HIS:HD2	1:A:761:HIS:CE1	1.91	0.88
1:A:123[A]:ARG:NH2	1:A:240:ASN:HB3	1.95	0.81
1:A:735:GLN:HE21	1:A:737:HIS:HB2	1.52	0.73
1:A:817:LYS:HZ3	7:A:1828:PEG:H22	1.56	0.70
1:A:569:ARG:HH22	6:A:1827:PGE:H12	1.54	0.68
1:A:692:HIS:CD2	1:A:761:HIS:CE1	2.69	0.63
1:A:731:HIS:HE1	7:A:1828:PEG:H31	1.65	0.62
1:A:123[A]:ARG:O	1:A:124:ASP:HB2	1.99	0.61
1:A:50:ILE:HD11	1:A:112:ILE:HD13	1.83	0.60
1:A:43:LEU:O	1:A:43:LEU:HG	2.04	0.58
1:A:43:LEU:HD22	1:A:94:LEU:HD23	1.86	0.58
1:A:690:GLU:OE2	1:A:692:HIS:HE1	1.90	0.54
1:A:193:PRO:HB2	1:A:205:PHE:HB3	1.90	0.53
1:A:643:ASP:OD1	1:A:761:HIS:HD2	1.91	0.53
1:A:569:ARG:NH2	6:A:1827:PGE:C1	2.63	0.53
1:A:731:HIS:HE1	7:A:1828:PEG:C3	2.21	0.52
1:A:643:ASP:CG	1:A:761:HIS:HD2	2.17	0.52
1:A:579:TYR:CZ	1:A:583:LEU:HD11	2.45	0.52
1:A:767:TYR:OH	7:A:1828:PEG:H42	2.12	0.49
1:A:767:TYR:OH	7:A:1828:PEG:C4	2.61	0.48
1:A:462:MET:O	1:A:476:PRO:HA	2.13	0.48
1:A:423:THR:HB	1:A:430:ALA:HB2	1.95	0.48
1:A:217:LYS:HE2	8:A:2075:HOH:O	2.13	0.47
1:A:511:GLY:H	1:A:527:TRP:CG	2.32	0.47
1:A:101:LEU:HB2	1:A:112:ILE:HB	1.96	0.47
1:A:230:GLY:H	4:A:1819:EDO:H21	1.80	0.47
1:A:555:GLU:HG3	8:A:2361:HOH:O	2.14	0.47
1:A:243:PRO:HB3	1:A:586:MET:HE1	1.97	0.47
1:A:93:GLN:O	1:A:93:GLN:HG3	2.15	0.46
1:A:643:ASP:OD1	1:A:761:HIS:CD2	2.69	0.46
1:A:123[A]:ARG:CZ	1:A:240:ASN:HB3	2.46	0.45
1:A:262:TRP:CG	1:A:337:LYS:HE3	2.52	0.45
1:A:569:ARG:NH1	8:A:2367:HOH:O	2.47	0.45
1:A:817:LYS:NZ	7:A:1828:PEG:H22	2.30	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:LEU:O	1:A:43:LEU:CG	2.66	0.44
1:A:507:HIS:CG	1:A:536:VAL:HB	2.53	0.44
1:A:365:ASP:HB2	1:A:366:PRO:HD2	2.00	0.43
1:A:468:GLY:O	1:A:471:ARG:HG2	2.20	0.42
1:A:534:GLN:HB2	1:A:535:PRO:CD	2.50	0.42
1:A:78:LEU:HD22	1:A:427:ILE:HD12	2.03	0.41
1:A:431:ASP:HB2	8:A:2281:HOH:O	2.20	0.41
1:A:332:LYS:HD3	1:A:332:LYS:HA	1.91	0.41
1:A:569:ARG:NH2	8:A:2371:HOH:O	2.54	0.41
1:A:258:LEU:O	1:A:457:ARG:HD3	2.20	0.40
1:A:450:LEU:HD23	1:A:450:LEU:HA	1.89	0.40
1:A:105:GLY:O	1:A:106:GLY:C	2.65	0.40
1:A:731:HIS:HE1	7:A:1828:PEG:C4	2.34	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	782/817 (96%)	755 (96%)	27 (4%)	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	643/690 (93%)	640 (100%)	3 (0%)	81	77

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	ASP
1	A	294	ASP
1	A	680	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	452	GLN
1	A	552	GLN
1	A	692	HIS
1	A	731	HIS
1	A	735	GLN
1	A	761	HIS
1	A	766	GLN
1	A	768	GLN
1	A	769	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](#)

12 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	EDO	A	1819	-	3,3,3	0.47	0	2,2,2	0.35	0
6	PGE	A	1827	-	9,9,9	0.51	0	8,8,8	0.24	0
3	5GF	A	1818	1	9,12,13	1.20	1 (11%)	14,18,20	1.84	4 (28%)
5	SO4	A	1820	-	4,4,4	0.24	0	6,6,6	0.09	0
5	SO4	A	1823	-	4,4,4	0.26	0	6,6,6	0.16	0
5	SO4	A	1822	-	4,4,4	0.25	0	6,6,6	0.08	0
5	SO4	A	1821	-	4,4,4	0.25	0	6,6,6	0.09	0
2	ARG	A	1829	-	10,11,11	0.77	1 (10%)	9,13,13	0.96	1 (11%)
5	SO4	A	1826	-	4,4,4	0.24	0	6,6,6	0.06	0
7	PEG	A	1828	-	6,6,6	0.69	0	5,5,5	0.49	0
5	SO4	A	1825	-	4,4,4	0.24	0	6,6,6	0.09	0
5	SO4	A	1824	-	4,4,4	0.25	0	6,6,6	0.10	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	1819	-	-	1/1/1/1	-
6	PGE	A	1827	-	-	4/7/7/7	-
3	5GF	A	1818	1	-	0/2/23/26	0/1/1/1
7	PEG	A	1828	-	-	2/4/4/4	-
2	ARG	A	1829	-	-	2/11/11/11	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1818	5GF	O5-C5	3.23	1.44	1.37
2	A	1829	ARG	OXT-C	-2.17	1.23	1.30

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1818	5GF	C1-O5-C5	3.89	120.41	113.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1818	5GF	O5-C1-C2	3.37	116.44	111.06
3	A	1818	5GF	C1-C2-C3	2.86	113.81	109.64
2	A	1829	ARG	OXT-C-O	-2.74	117.87	124.08
3	A	1818	5GF	C5-C4-C3	-2.57	107.88	110.71

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1827	PGE	O2-C3-C4-O3
6	A	1827	PGE	O1-C1-C2-O2
7	A	1828	PEG	O1-C1-C2-O2
2	A	1829	ARG	OXT-C-CA-N
6	A	1827	PGE	O3-C5-C6-O4
6	A	1827	PGE	C1-C2-O2-C3
7	A	1828	PEG	C4-C3-O2-C2
4	A	1819	EDO	O1-C1-C2-O2
2	A	1829	ARG	O-C-CA-N

There are no ring outliers.

3 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1819	EDO	1	0
6	A	1827	PGE	3	0
7	A	1828	PEG	7	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	781/817 (95%)	1.29	153 (19%) 3 2	14, 28, 54, 101	5 (0%)

All (153) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	137	VAL	10.2
1	A	376	TYR	7.6
1	A	43	LEU	6.7
1	A	138	ASN	5.9
1	A	101	LEU	5.6
1	A	35	ALA	5.6
1	A	106	GLY	5.1
1	A	95	THR	5.1
1	A	89	ALA	5.0
1	A	90	ILE	4.9
1	A	105	GLY	4.8
1	A	44	ARG	4.8
1	A	97	THR	4.7
1	A	48	LEU	4.7
1	A	94	LEU	4.6
1	A	45	ASN	4.5
1	A	37	PHE	4.4
1	A	117	PHE	4.4
1	A	367	GLN	4.4
1	A	49	VAL	4.3
1	A	701	LEU	4.2
1	A	785	ALA	4.1
1	A	47	GLY	4.0
1	A	91	LYS	4.0
1	A	59	LEU	4.0
1	A	112	ILE	4.0
1	A	54	GLN	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	46	ASP	3.9
1	A	319	GLU	3.8
1	A	42	LYS	3.8
1	A	142	ILE	3.7
1	A	74	ASP	3.7
1	A	93	GLN	3.7
1	A	76	LYS	3.7
1	A	69	VAL	3.6
1	A	85	GLN	3.6
1	A	87	GLU	3.6
1	A	52	THR	3.6
1	A	125	SER	3.5
1	A	116	PRO	3.5
1	A	236	LEU	3.5
1	A	816	HIS	3.5
1	A	38	TYR	3.5
1	A	92	ALA	3.5
1	A	124	ASP	3.4
1	A	784[A]	CYS	3.4
1	A	40	GLU	3.3
1	A	56	ASN	3.3
1	A	777	ILE	3.3
1	A	769	HIS	3.3
1	A	104	SER	3.3
1	A	36	VAL	3.2
1	A	107	THR	3.2
1	A	50	ILE	3.2
1	A	98	GLU	3.2
1	A	375	LEU	3.2
1	A	748	ARG	3.1
1	A	39	LYS	3.1
1	A	240	ASN	3.1
1	A	86	PRO	3.1
1	A	790	LEU	3.1
1	A	61	PHE	3.1
1	A	136	GLN	3.1
1	A	109	THR	3.1
1	A	771	ASP	3.1
1	A	377	PHE	3.0
1	A	776	THR	3.0
1	A	368	GLY	3.0
1	A	130	ALA	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	115	ARG	3.0
1	A	447	GLN	3.0
1	A	123[A]	ARG	3.0
1	A	118	ALA	3.0
1	A	102	GLN	2.9
1	A	53	ASN	2.9
1	A	57	ILE	2.9
1	A	111	ARG	2.9
1	A	127	LEU	2.9
1	A	787	THR	2.8
1	A	649	ASP	2.8
1	A	768	GLN	2.7
1	A	51	THR	2.7
1	A	812	ASN	2.7
1	A	789	ALA	2.7
1	A	103	PHE	2.7
1	A	150	PRO	2.7
1	A	808	ARG	2.7
1	A	143	ASN	2.7
1	A	374	GLU	2.7
1	A	775	LYS	2.7
1	A	65	ALA	2.6
1	A	601	ALA	2.6
1	A	778	ALA	2.6
1	A	122	TYR	2.6
1	A	228	LYS	2.6
1	A	220	SER	2.6
1	A	41	HIS	2.6
1	A	70	LEU	2.6
1	A	113	GLN	2.6
1	A	96	GLU	2.6
1	A	64	GLU	2.5
1	A	660	HIS	2.5
1	A	372	ALA	2.5
1	A	135	PHE	2.4
1	A	737	HIS	2.4
1	A	333	GLN	2.4
1	A	141	LYS	2.4
1	A	72	ARG	2.4
1	A	71	TYR	2.4
1	A	120	SER	2.4
1	A	114	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	128	LEU	2.3
1	A	110	ALA	2.3
1	A	366	PRO	2.3
1	A	774	GLY	2.3
1	A	643	ASP	2.3
1	A	450	LEU	2.3
1	A	99	ASN	2.3
1	A	421	GLU	2.3
1	A	100	HIS	2.3
1	A	121	TYR	2.3
1	A	378	GLY	2.3
1	A	770	LEU	2.3
1	A	75	SER	2.2
1	A	733	ASP	2.2
1	A	83	LEU	2.2
1	A	451	ASP	2.2
1	A	119	ILE	2.2
1	A	634	LYS	2.2
1	A	647	GLN	2.2
1	A	88	SER	2.2
1	A	370	PRO	2.2
1	A	600	PRO	2.2
1	A	685	ARG	2.2
1	A	151	GLY	2.2
1	A	82	ALA	2.2
1	A	147	TYR	2.1
1	A	772	ILE	2.1
1	A	400	ASP	2.1
1	A	63	SER	2.1
1	A	149	SER	2.1
1	A	126	GLU	2.1
1	A	365	ASP	2.1
1	A	79	PRO	2.1
1	A	108	LEU	2.1
1	A	60	GLN	2.1
1	A	369	GLN	2.1
1	A	721	GLN	2.1
1	A	735	GLN	2.1
1	A	659	LEU	2.1
1	A	703	GLN	2.0
1	A	628	VAL	2.0
1	A	813	LEU	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($\text{\AA}^2$)	Q<0.9
5	SO4	A	1826	5/5	0.62	0.17	95,98,99,101	0
5	SO4	A	1821	5/5	0.66	0.19	78,80,85,90	0
4	EDO	A	1819	4/4	0.69	0.27	43,53,59,59	0
5	SO4	A	1825	5/5	0.72	0.17	70,74,80,83	0
2	ARG	A	1829	12/12	0.72	0.17	27,51,57,59	0
7	PEG	A	1828	7/7	0.76	0.21	25,32,39,43	0
5	SO4	A	1820	5/5	0.81	0.14	67,68,74,77	0
6	PGE	A	1827	10/10	0.84	0.22	31,49,57,57	0
5	SO4	A	1824	5/5	0.85	0.14	72,73,79,80	0
5	SO4	A	1823	5/5	0.91	0.11	51,52,59,62	0
3	5GF	A	1818	12/13	0.91	0.10	17,20,23,23	0
5	SO4	A	1822	5/5	0.96	0.10	29,31,38,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.