



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:13 PM UTC

PDB ID : 6ZJO / pdb_00006zjo
Title : Crystal Structure of Staphylococcus aureus RsgA.
Authors : Bennison, D.J.; Rafferty, J.B.; Corrigan, R.M.
Deposited on : 2020-06-29
Resolution : 2.01 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

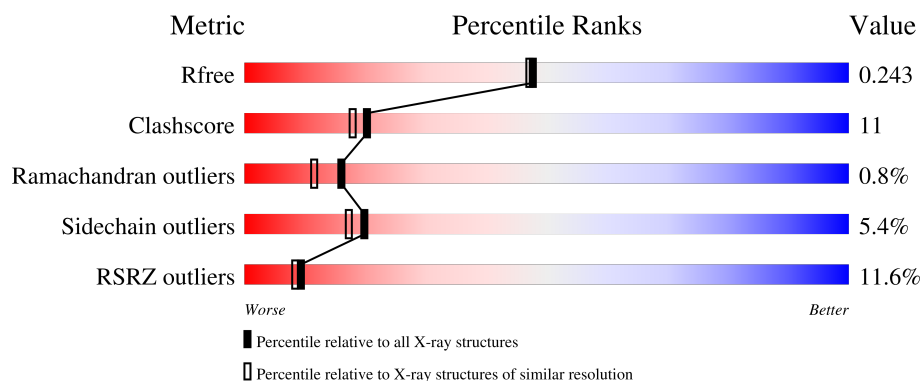
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	311	9% 68% 14% •• 15%
1	B	311	11% 67% 17% • 14%

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
4	PO4	B	311	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4564 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 Small ribosomal subunit biogenesis GTPase RsgA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	0	0
			2112	1347	356	404	5			
1	B	268	Total	C	N	O	S	0	0	0
			2159	1377	375	402	5			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A0A4V5LHH4
A	-18	GLY	-	expression tag	UNP A0A4V5LHH4
A	-17	SER	-	expression tag	UNP A0A4V5LHH4
A	-16	SER	-	expression tag	UNP A0A4V5LHH4
A	-15	HIS	-	expression tag	UNP A0A4V5LHH4
A	-14	HIS	-	expression tag	UNP A0A4V5LHH4
A	-13	HIS	-	expression tag	UNP A0A4V5LHH4
A	-12	HIS	-	expression tag	UNP A0A4V5LHH4
A	-11	HIS	-	expression tag	UNP A0A4V5LHH4
A	-10	HIS	-	expression tag	UNP A0A4V5LHH4
A	-9	SER	-	expression tag	UNP A0A4V5LHH4
A	-8	SER	-	expression tag	UNP A0A4V5LHH4
A	-7	GLY	-	expression tag	UNP A0A4V5LHH4
A	-6	LEU	-	expression tag	UNP A0A4V5LHH4
A	-5	VAL	-	expression tag	UNP A0A4V5LHH4
A	-4	PRO	-	expression tag	UNP A0A4V5LHH4
A	-3	ARG	-	expression tag	UNP A0A4V5LHH4
A	-2	GLY	-	expression tag	UNP A0A4V5LHH4
A	-1	SER	-	expression tag	UNP A0A4V5LHH4
A	0	HIS	-	expression tag	UNP A0A4V5LHH4
B	-19	MET	-	initiating methionine	UNP A0A4V5LHH4
B	-18	GLY	-	expression tag	UNP A0A4V5LHH4
B	-17	SER	-	expression tag	UNP A0A4V5LHH4
B	-16	SER	-	expression tag	UNP A0A4V5LHH4
B	-15	HIS	-	expression tag	UNP A0A4V5LHH4

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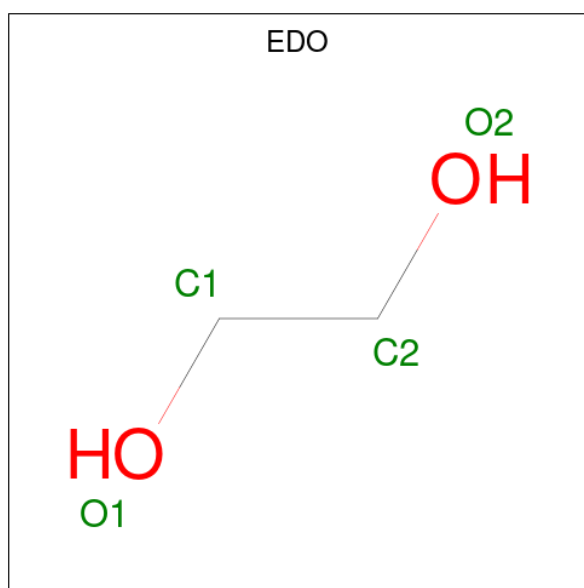
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP A0A4V5LHH4
B	-13	HIS	-	expression tag	UNP A0A4V5LHH4
B	-12	HIS	-	expression tag	UNP A0A4V5LHH4
B	-11	HIS	-	expression tag	UNP A0A4V5LHH4
B	-10	HIS	-	expression tag	UNP A0A4V5LHH4
B	-9	SER	-	expression tag	UNP A0A4V5LHH4
B	-8	SER	-	expression tag	UNP A0A4V5LHH4
B	-7	GLY	-	expression tag	UNP A0A4V5LHH4
B	-6	LEU	-	expression tag	UNP A0A4V5LHH4
B	-5	VAL	-	expression tag	UNP A0A4V5LHH4
B	-4	PRO	-	expression tag	UNP A0A4V5LHH4
B	-3	ARG	-	expression tag	UNP A0A4V5LHH4
B	-2	GLY	-	expression tag	UNP A0A4V5LHH4
B	-1	SER	-	expression tag	UNP A0A4V5LHH4
B	0	HIS	-	expression tag	UNP A0A4V5LHH4

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

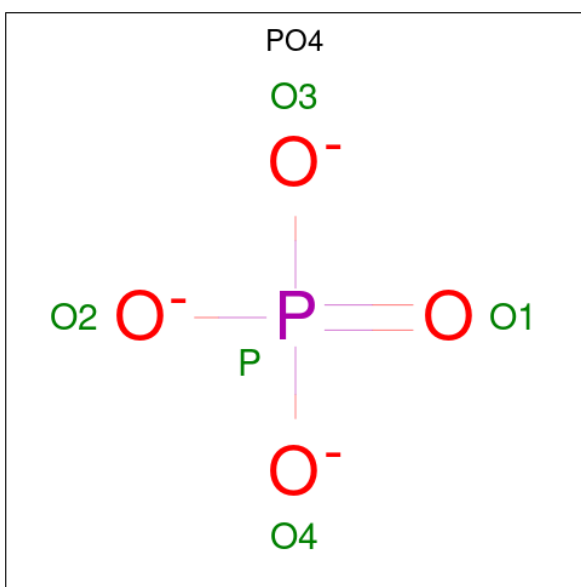
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

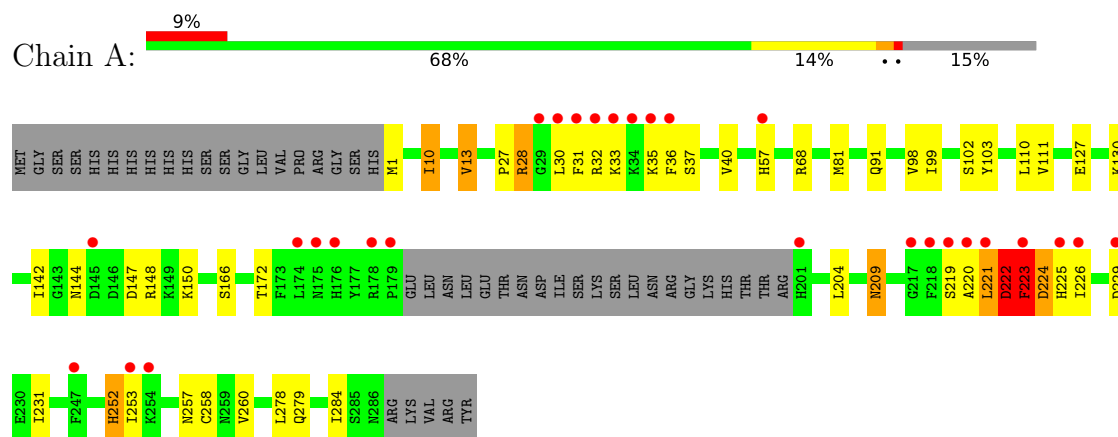
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	102	Total	O	0	0
			102	102		
5	B	110	Total	O	0	0
			110	110		

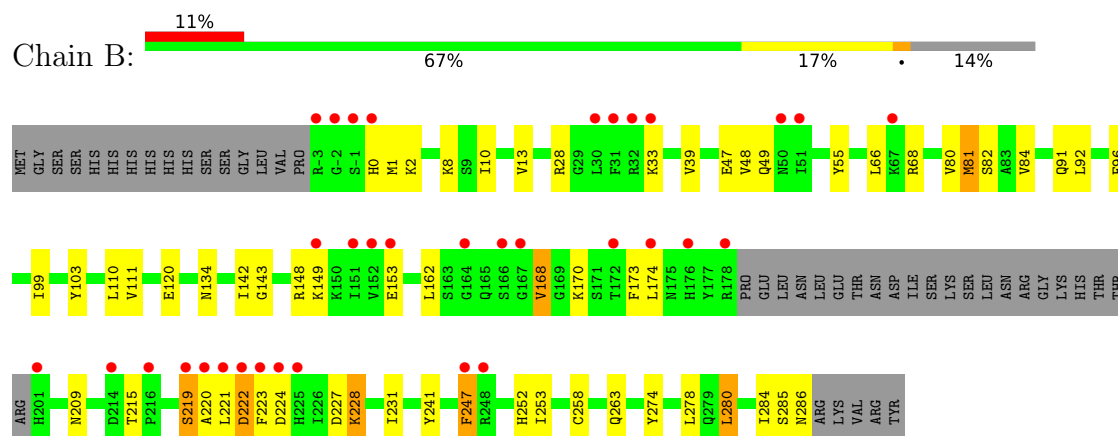
3 Residue-property plots [i](#)

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: Small ribosomal subunit biogenesis GTPase RsgA



- Molecule 1: Small ribosomal subunit biogenesis GTPase RsgA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.67Å 93.57Å 68.18Å 90.00° 90.71° 90.00°	Depositor
Resolution (Å)	47.20 – 2.01 47.20 – 2.01	Depositor EDS
% Data completeness (in resolution range)	98.2 (47.20-2.01) 98.0 (47.20-2.01)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.224 , 0.278 (Not available) , 0.243	Depositor DCC
R_{free} test set	2186 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.140	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 32.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.106 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4564	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.46% 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: PO4, ZN, 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	1.03	0/2157	1.36	6/2925 (0.2%)
1	B	1.06	0/2203	1.38	4/2975 (0.1%)
All	All	1.05	0/4360	1.37	10/5900 (0.2%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	223	PHE	N-CA-C	9.20	122.29	111.71
1	B	223	PHE	N-CA-C	-6.44	101.73	110.68
1	A	28	ARG	CA-C-N	5.78	126.50	120.03
1	A	28	ARG	C-N-CA	5.78	126.50	120.03
1	A	222	ASP	N-CA-C	5.69	118.82	109.72
1	A	221	LEU	N-CA-C	5.21	117.44	110.35
1	A	224	ASP	N-CA-C	-5.13	99.88	110.80
1	B	66	LEU	CA-C-O	-5.07	115.43	121.16
1	B	143	GLY	CA-C-N	5.03	127.27	120.38
1	B	143	GLY	C-N-CA	5.03	127.27	120.38

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	2112	0	2007	51	0
1	B	2159	0	2100	45	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	28	0	42	0	0
3	B	36	0	54	2	0
4	A	5	0	0	0	0
4	B	10	0	0	3	0
5	A	102	0	0	4	0
5	B	110	0	0	4	0
All	All	4564	0	4203	97	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 11.

All (97) 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:98:VAL:HG12	1:A:223:PHE:CZ	1.62	1.34
1:A:98:VAL:HG12	1:A:223:PHE:HZ	0.94	1.08
1:A:98:VAL:CG1	1:A:223:PHE:HZ	1.64	1.08
1:A:31:PHE:HB3	1:A:36:PHE:CD1	1.92	1.05
1:A:253:ILE:HD13	1:A:278:LEU:CD1	1.95	0.97
1:A:253:ILE:CD1	1:A:278:LEU:CD1	2.43	0.96
1:A:31:PHE:HB3	1:A:36:PHE:CE1	2.00	0.94
1:A:102:SER:HB3	1:A:223:PHE:HE1	1.34	0.92
1:A:98:VAL:CG1	1:A:223:PHE:CZ	2.47	0.92
1:B:162:LEU:O	1:B:215:THR:HG22	1.77	0.84
1:B:285:SER:O	1:B:286:ASN:OD1	1.96	0.82
1:A:253:ILE:HD13	1:A:278:LEU:HD13	1.63	0.79
1:B:134:ASN:HB3	5:B:481:HOH:O	1.83	0.78
1:A:31:PHE:CB	1:A:36:PHE:CE1	2.69	0.76
1:A:102:SER:HB3	1:A:223:PHE:CE1	2.22	0.74
1:A:253:ILE:CD1	1:A:278:LEU:HD12	2.22	0.68
1:B:10:ILE:O	1:B:13:VAL:HG12	1.95	0.67
1:A:81:MET:O	1:A:111:VAL:HA	1.96	0.65
1:B:241:TYR:CD1	1:B:263:GLN:HG3	2.32	0.64
1:A:253:ILE:CD1	1:A:278:LEU:HD11	2.28	0.62
1:B:170:LYS:CB	4:B:311:PO4:O3	2.48	0.62
1:B:103:TYR:OH	1:B:221:LEU:HD22	2.00	0.61
1:B:1:MET:HE2	1:B:47:GLU:HB2	1.86	0.58
1:B:103:TYR:OH	1:B:221:LEU:CD2	2.52	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:253:ILE:HD12	1:A:278:LEU:CD1	2.29	0.57
1:A:231:ILE:HD11	1:A:284:ILE:HG13	1.87	0.57
1:A:35:LYS:CB	5:A:485:HOH:O	2.53	0.57
1:A:27:PRO:HB3	1:A:36:PHE:CE1	2.40	0.56
1:B:68:ARG:NH2	1:B:221:LEU:HD23	2.20	0.56
1:B:81:MET:O	1:B:111:VAL:HA	2.06	0.56
1:A:102:SER:CB	1:A:223:PHE:HE1	2.13	0.55
1:B:170:LYS:HB3	4:B:311:PO4:O3	2.06	0.55
1:B:247:PHE:HB2	5:B:429:HOH:O	2.06	0.55
1:A:102:SER:CB	1:A:223:PHE:CE1	2.89	0.54
1:A:103:TYR:OH	1:A:224:ASP:HB2	2.08	0.54
1:B:2:LYS:HE2	1:B:48:VAL:HG21	1.90	0.53
1:A:221:LEU:HD12	1:A:222:ASP:H	1.72	0.53
1:A:32:ARG:HA	1:A:36:PHE:CD2	2.42	0.53
1:B:0:HIS:NE2	1:B:47:GLU:OE2	2.42	0.53
1:B:99:ILE:HD11	1:B:222:ASP:CG	2.34	0.53
1:B:80:VAL:HG12	1:B:168:VAL:CG1	2.39	0.52
1:B:221:LEU:HA	1:B:224:ASP:CB	2.40	0.52
1:B:162:LEU:O	1:B:215:THR:CG2	2.55	0.52
1:A:91:GLN:NE2	1:A:279:GLN:OE1	2.42	0.51
1:B:110:LEU:HD21	1:B:173:PHE:CZ	2.46	0.51
1:A:219:SER:O	1:A:221:LEU:N	2.45	0.50
1:B:280:LEU:O	1:B:284:ILE:HG12	2.11	0.50
1:B:231:ILE:HD11	1:B:284:ILE:HG13	1.94	0.50
1:A:147:ASP:OD2	1:A:150:LYS:CB	2.60	0.49
1:B:219:SER:H	1:B:221:LEU:HD12	1.76	0.49
1:A:253:ILE:HD13	1:A:278:LEU:HD11	1.85	0.48
1:B:219:SER:C	1:B:221:LEU:H	2.21	0.48
1:A:127:GLU:OE1	1:A:130:LYS:NZ	2.40	0.48
1:A:68:ARG:CB	5:A:493:HOH:O	2.62	0.48
1:A:27:PRO:HB3	1:A:36:PHE:HE1	1.78	0.48
1:A:33:LYS:N	1:A:36:PHE:HB3	2.29	0.47
1:B:148:ARG:O	1:B:149:LYS:C	2.58	0.47
1:A:253:ILE:HD12	1:A:278:LEU:HD11	1.95	0.47
1:B:81:MET:HE3	1:B:81:MET:HB3	1.77	0.47
1:A:148:ARG:HD2	1:A:172:THR:HG22	1.95	0.47
1:B:170:LYS:N	4:B:311:PO4:O3	2.38	0.47
1:B:220:ALA:O	1:B:224:ASP:CB	2.64	0.46
1:A:28:ARG:HD2	1:A:57:HIS:CE1	2.51	0.46
1:B:228:LYS:HE2	1:B:285:SER:OG	2.17	0.45
1:B:120:GLU:HB2	3:B:302:EDO:H22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:429:HOH:O	3:B:302:EDO:H21	2.16	0.44
1:B:80:VAL:HG12	1:B:168:VAL:HG12	1.98	0.44
1:A:252:HIS:HB3	1:A:260:VAL:HB	1.98	0.44
1:A:33:LYS:H	1:A:36:PHE:HB3	1.82	0.44
1:A:209:ASN:ND2	5:A:410:HOH:O	2.50	0.44
1:B:274:TYR:CZ	1:B:278:LEU:HD13	2.53	0.44
1:A:103:TYR:HE2	1:A:222:ASP:O	2.00	0.44
1:B:92:LEU:HG	1:B:96:PHE:CE2	2.53	0.44
1:B:142:ILE:HD11	1:B:173:PHE:HZ	1.82	0.44
1:B:228:LYS:HG2	5:B:467:HOH:O	2.16	0.44
1:A:1:MET:HE1	1:A:57:HIS:CG	2.53	0.43
1:B:285:SER:C	1:B:286:ASN:OD1	2.59	0.43
1:B:153:GLU:HA	5:B:402:HOH:O	2.18	0.43
1:A:103:TYR:CE2	1:A:222:ASP:O	2.71	0.43
1:B:241:TYR:CG	1:B:263:GLN:HG3	2.53	0.43
1:A:36:PHE:CD2	1:A:36:PHE:O	2.73	0.42
1:A:102:SER:HG	1:A:223:PHE:HD1	1.63	0.42
1:A:110:LEU:HD11	1:A:142:ILE:HG12	2.02	0.42
1:B:8:LYS:HD2	1:B:10:ILE:CD1	2.49	0.42
1:A:81:MET:HE3	1:A:81:MET:HB3	1.83	0.42
1:A:252:HIS:HB3	1:A:258:CYS:SG	2.60	0.42
1:B:222:ASP:C	1:B:224:ASP:H	2.28	0.42
1:A:31:PHE:O	1:A:36:PHE:CZ	2.73	0.42
1:B:110:LEU:HD11	1:B:142:ILE:HG12	2.02	0.41
1:A:99:ILE:HA	1:A:223:PHE:CE1	2.56	0.41
1:B:49:GLN:HG3	1:B:55:TYR:CD2	2.56	0.41
1:A:10:ILE:O	1:A:13:VAL:CG1	2.69	0.41
1:A:10:ILE:O	1:A:13:VAL:HG12	2.21	0.41
1:B:82:SER:HB2	1:B:84:VAL:O	2.21	0.40
1:B:28:ARG:HA	1:B:55:TYR:CE1	2.57	0.40
1:A:40:VAL:HG13	1:A:204:LEU:HG	2.04	0.40
1:B:252:HIS:HA	1:B:258:CYS:SG	2.62	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	261/311 (84%)	244 (94%)	15 (6%)	2 (1%)	16	11
1	B	264/311 (85%)	243 (92%)	19 (7%)	2 (1%)	16	11
All	All	525/622 (84%)	487 (93%)	34 (6%)	4 (1%)	16	11

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	219	SER
1	A	220	ALA
1	B	33	LYS
1	A	10	ILE

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	227/283 (80%)	214 (94%)	13 (6%)	18	15
1	B	232/283 (82%)	220 (95%)	12 (5%)	21	18
All	All	459/566 (81%)	434 (95%)	25 (5%)	20	17

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	VAL
1	A	30	LEU
1	A	37	SER
1	A	144	ASN
1	A	166	SER
1	A	209	ASN
1	A	222	ASP
1	A	223	PHE

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Mol	Chain	Res	Type
1	A	225	HIS
1	A	226	ILE
1	A	229	ASP
1	A	252	HIS
1	A	257	ASN
1	B	39	VAL
1	B	81	MET
1	B	91	GLN
1	B	168	VAL
1	B	174	LEU
1	B	209	ASN
1	B	222	ASP
1	B	227	ASP
1	B	228	LYS
1	B	247	PHE
1	B	253	ILE
1	B	280	LEU

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

Mol	Chain	Res	Type
1	A	57	HIS
1	B	49	GLN
1	B	91	GLN
1	B	144	ASN
1	B	271	GLN
1	B	279	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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 21 ligands modelled in this entry, 2 are monoatomic - leaving 19 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	A	308	-	3,3,3	0.13	0	2,2,2	0.09	0
3	EDO	A	305	-	3,3,3	0.27	0	2,2,2	0.24	0
3	EDO	B	310	-	3,3,3	0.40	0	2,2,2	0.51	0
3	EDO	A	304	-	3,3,3	0.56	0	2,2,2	0.64	0
3	EDO	A	302	-	3,3,3	0.11	0	2,2,2	0.42	0
3	EDO	B	307	-	3,3,3	0.13	0	2,2,2	0.21	0
3	EDO	B	303	-	3,3,3	0.08	0	2,2,2	0.12	0
3	EDO	B	308	-	3,3,3	0.09	0	2,2,2	0.14	0
3	EDO	B	304	-	3,3,3	0.09	0	2,2,2	0.16	0
4	PO4	B	311	-	4,4,4	0.96	0	6,6,6	0.46	0
3	EDO	B	302	-	3,3,3	0.55	0	2,2,2	0.81	0
4	PO4	B	312	-	4,4,4	0.54	0	6,6,6	0.52	0
3	EDO	A	307	-	3,3,3	0.14	0	2,2,2	0.31	0
3	EDO	B	305	-	3,3,3	0.12	0	2,2,2	0.16	0
3	EDO	A	303	-	3,3,3	0.15	0	2,2,2	0.27	0
3	EDO	B	306	-	3,3,3	0.13	0	2,2,2	0.11	0
3	EDO	B	309	-	3,3,3	0.20	0	2,2,2	0.11	0
4	PO4	A	309	-	4,4,4	0.92	0	6,6,6	0.82	0
3	EDO	A	306	-	3,3,3	0.17	0	2,2,2	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
3	EDO	A	308	-	-	1/1/1/1	-
3	EDO	B	307	-	-	1/1/1/1	-
3	EDO	A	305	-	-	1/1/1/1	-
3	EDO	A	307	-	-	1/1/1/1	-
3	EDO	B	305	-	-	0/1/1/1	-
3	EDO	B	302	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	B	310	-	-	1/1/1/1	-
3	EDO	A	303	-	-	1/1/1/1	-
3	EDO	A	304	-	-	0/1/1/1	-
3	EDO	B	303	-	-	1/1/1/1	-
3	EDO	B	304	-	-	1/1/1/1	-
3	EDO	B	306	-	-	1/1/1/1	-
3	EDO	B	308	-	-	0/1/1/1	-
3	EDO	B	309	-	-	0/1/1/1	-
3	EDO	A	302	-	-	1/1/1/1	-
3	EDO	A	306	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	304	EDO	O1-C1-C2-O2
3	A	303	EDO	O1-C1-C2-O2
3	A	308	EDO	O1-C1-C2-O2
3	B	307	EDO	O1-C1-C2-O2
3	A	305	EDO	O1-C1-C2-O2
3	A	306	EDO	O1-C1-C2-O2
3	B	303	EDO	O1-C1-C2-O2
3	A	302	EDO	O1-C1-C2-O2
3	B	306	EDO	O1-C1-C2-O2
3	A	307	EDO	O1-C1-C2-O2
3	B	310	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	311	PO4	3	0
3	B	302	EDO	2	0

5.7 Other polymers [i](#)

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	265/311 (85%)	0.54	28 (10%)	11 10	22, 35, 74, 113	0
1	B	268/311 (86%)	0.62	34 (12%)	8 7	19, 36, 70, 94	0
All	All	533/622 (85%)	0.58	62 (11%)	9 8	19, 35, 71, 113	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	223	PHE	10.9
1	A	36	PHE	9.3
1	A	219	SER	7.9
1	A	218	PHE	6.8
1	B	31	PHE	6.1
1	A	221	LEU	5.9
1	B	247	PHE	5.5
1	B	223	PHE	5.3
1	B	224	ASP	5.2
1	B	30	LEU	5.2
1	A	31	PHE	5.1
1	A	253	ILE	5.0
1	A	225	HIS	4.8
1	A	32	ARG	4.5
1	A	220	ALA	4.4
1	B	222	ASP	4.3
1	B	51	ILE	4.3
1	A	34	LYS	4.3
1	A	35	LYS	4.1
1	A	254	LYS	4.1
1	B	220	ALA	4.0
1	B	225	HIS	3.8
1	A	179	PRO	3.6
1	B	216	PRO	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	33	LYS	3.5
1	B	201	HIS	3.4
1	A	145	ASP	3.3
1	A	201	HIS	3.1
1	A	30	LEU	2.9
1	B	214	ASP	2.8
1	A	174	LEU	2.8
1	B	50	ASN	2.7
1	A	217	GLY	2.7
1	A	226	ILE	2.7
1	B	152	VAL	2.7
1	B	33	LYS	2.7
1	B	149	LYS	2.6
1	B	219	SER	2.6
1	B	-3	ARG	2.6
1	B	178	ARG	2.6
1	B	0	HIS	2.6
1	A	29	GLY	2.5
1	B	174	LEU	2.5
1	B	32	ARG	2.5
1	B	172	THR	2.5
1	B	-2	GLY	2.5
1	A	229	ASP	2.3
1	B	166	SER	2.3
1	B	221	LEU	2.3
1	B	151	ILE	2.3
1	B	248	ARG	2.3
1	B	167	GLY	2.3
1	A	247	PHE	2.2
1	A	176	HIS	2.2
1	B	164	GLY	2.2
1	A	175	ASN	2.1
1	B	-1	SER	2.1
1	A	178	ARG	2.0
1	A	57	HIS	2.0
1	B	67	LYS	2.0
1	B	153	GLU	2.0
1	B	176	HIS	2.0

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

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
3	EDO	B	307	4/4	0.57	0.20	87,88,90,94	0
3	EDO	A	307	4/4	0.64	0.18	65,77,78,78	0
3	EDO	B	305	4/4	0.68	0.17	62,67,78,84	0
3	EDO	A	303	4/4	0.68	0.22	90,91,96,98	0
3	EDO	B	306	4/4	0.70	0.22	60,72,74,78	0
3	EDO	B	304	4/4	0.70	0.20	54,55,57,58	0
3	EDO	A	308	4/4	0.78	0.17	61,63,66,67	0
3	EDO	B	308	4/4	0.78	0.17	67,67,68,70	0
4	PO4	B	311	5/5	0.79	0.12	45,65,67,68	0
3	EDO	B	310	4/4	0.80	0.16	49,51,51,53	0
3	EDO	A	302	4/4	0.81	0.16	46,47,50,52	0
3	EDO	B	303	4/4	0.85	0.17	49,56,57,59	0
3	EDO	B	302	4/4	0.85	0.13	35,35,35,41	0
3	EDO	A	306	4/4	0.87	0.12	52,57,61,62	0
3	EDO	B	309	4/4	0.87	0.14	46,46,47,49	0
3	EDO	A	304	4/4	0.87	0.14	43,44,44,46	0
3	EDO	A	305	4/4	0.87	0.17	42,48,53,58	0
4	PO4	B	312	5/5	0.88	0.12	50,50,53,63	0
4	PO4	A	309	5/5	0.98	0.04	33,33,39,39	0
2	ZN	B	301	1/1	0.99	0.02	34,34,34,34	0
2	ZN	A	301	1/1	0.99	0.04	35,35,35,35	0

6.5 Other polymers [i](#)

There are no such residues in this entry.