



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 12:29 pm GMT

PDB ID : 3CZP
Title : Crystal structure of putative polyphosphate kinase 2 from *Pseudomonas aeruginosa* PA01
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Deposited on : 2008-04-29
Resolution : 2.00 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

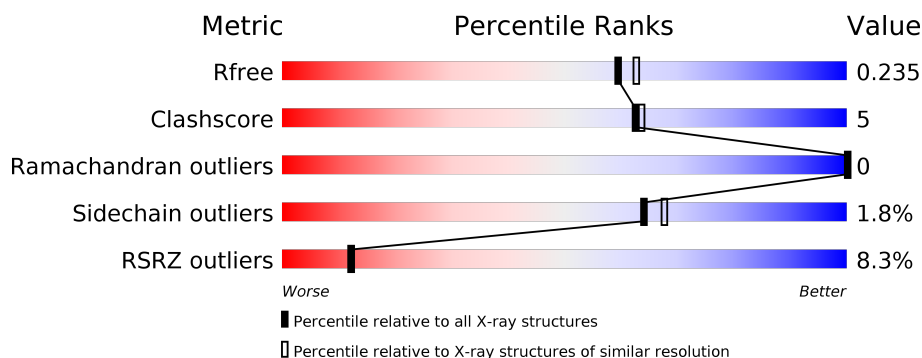
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (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 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	500	<div> <div>7%</div> <div> <div></div> <div>82%</div> <div>11%</div> <div>7%</div> </div> </div>
1	B	500	<div> <div>8%</div> <div> <div></div> <div>80%</div> <div>11%</div> <div>8%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	A	608	-	-	-	X
2	ACT	B	602	-	-	X	X
5	GOL	A	602	-	-	-	X
5	GOL	A	603	-	-	-	X
5	GOL	A	609	-	-	X	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8481 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.

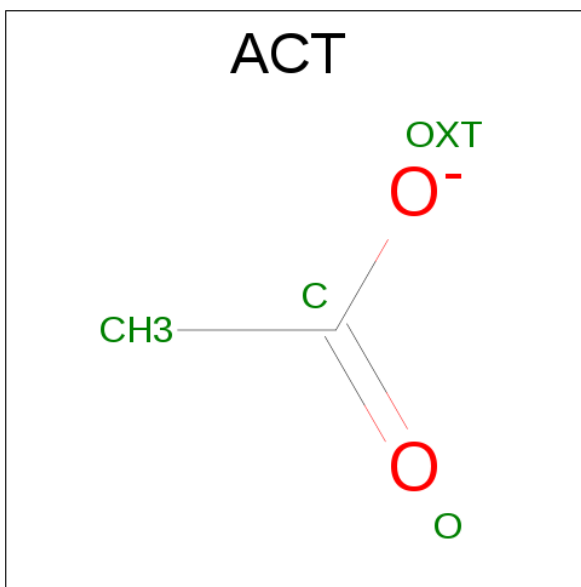
- Molecule 1 is a protein called Putative polyphosphate kinase 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	466	Total	C	N	O	S	Se	0	7	0
			3925	2507	701	708	2	7			
1	B	458	Total	C	N	O	S	Se	0	2	0
			3825	2446	684	686	2	7			

There are 8 discrepancies between the modelled and reference sequences:

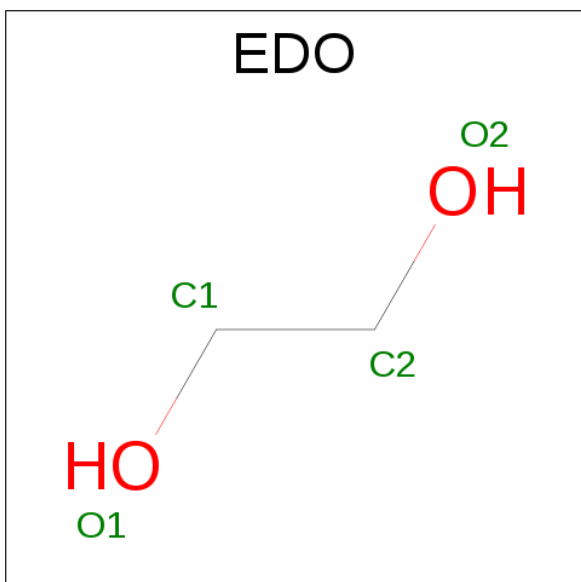
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP Q9HYF1
A	0	HIS	-	EXPRESSION TAG	UNP Q9HYF1
A	497	GLY	-	EXPRESSION TAG	UNP Q9HYF1
A	498	ALA	-	EXPRESSION TAG	UNP Q9HYF1
B	-1	GLY	-	EXPRESSION TAG	UNP Q9HYF1
B	0	HIS	-	EXPRESSION TAG	UNP Q9HYF1
B	497	GLY	-	EXPRESSION TAG	UNP Q9HYF1
B	498	ALA	-	EXPRESSION TAG	UNP Q9HYF1

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



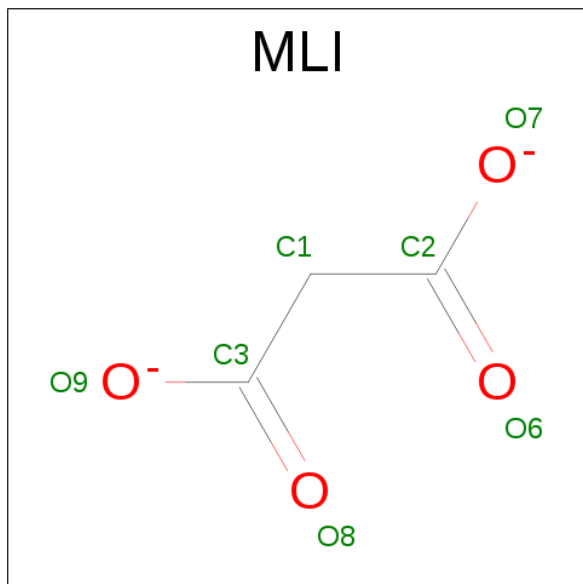
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		

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



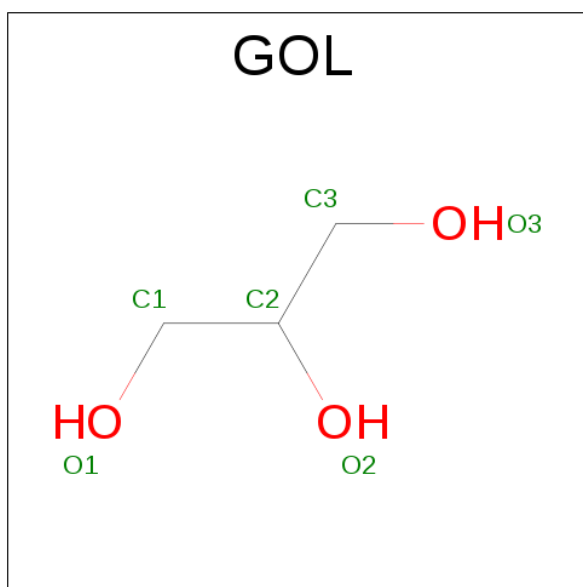
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is MALONATE ION (three-letter code: MLI) (formula: $C_3H_2O_4$).



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

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

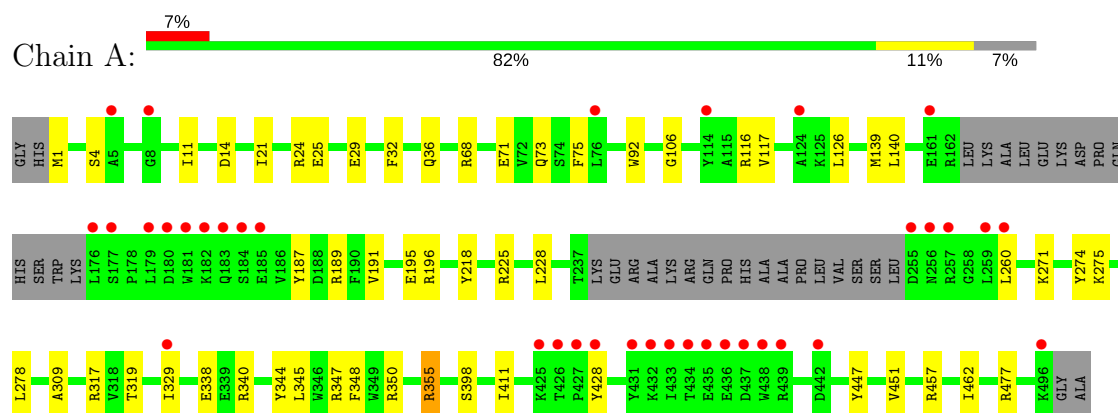
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	323	Total	O	0	0
			323	323		
6	B	345	Total	O	0	0
			345	345		

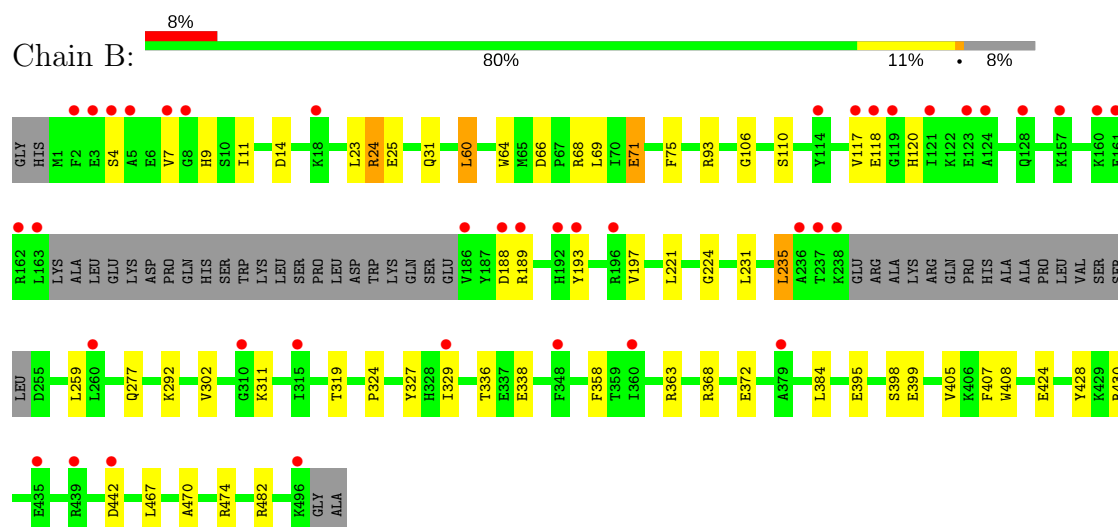
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Putative polyphosphate kinase 2



• Molecule 1: Putative polyphosphate kinase 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	97.09Å 100.73Å 120.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.00 38.59 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.0 (40.00-2.00) 94.7 (38.59-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.68 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.173 , 0.231 0.180 , 0.235	Depositor DCC
R_{free} test set	3811 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	31.8	Xtriage
Anisotropy	0.695	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.017 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8481	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% 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: GOL, MLI, EDO, 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	0.91	0/4027	0.83	2/5423 (0.0%)
1	B	0.87	4/3915 (0.1%)	0.83	3/5273 (0.1%)
All	All	0.89	4/7942 (0.1%)	0.83	5/10696 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	71	GLU	CG-CD	5.30	1.59	1.51
1	B	235	LEU	CB-CG	5.28	1.67	1.52
1	B	302	VAL	CB-CG1	5.17	1.63	1.52
1	B	470	ALA	CA-CB	5.13	1.63	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	93	ARG	NE-CZ-NH1	5.85	123.22	120.30
1	A	340	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	B	14	ASP	CB-CG-OD1	5.13	122.92	118.30
1	A	355	ARG	NE-CZ-NH2	-5.04	117.78	120.30
1	B	474	ARG	NE-CZ-NH2	-5.02	117.79	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	3925	0	3874	47	0
1	B	3825	0	3784	38	0
2	A	16	0	12	1	0
2	B	4	0	3	2	0
3	A	4	0	6	0	0
3	B	8	0	12	2	0
4	B	7	0	2	0	0
5	A	24	0	32	12	0
6	A	323	0	0	3	0
6	B	345	0	0	4	0
All	All	8481	0	7725	78	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 5.

All (78) 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:31:GLN:HE22	1:B:66:ASP:H	1.11	0.95
1:B:259:LEU:HD21	1:B:467[A]:LEU:HD11	1.58	0.86
1:A:225:ARG:HG2	5:A:602:GOL:H31	1.63	0.81
1:A:24:ARG:HD3	5:A:609:GOL:H11	1.67	0.76
1:B:277:GLN:HE21	1:B:482:ARG:HH12	1.37	0.71
1:B:424:GLU:HG2	6:B:947:HOH:O	1.90	0.71
1:B:259:LEU:CD2	1:B:467[A]:LEU:HD11	2.24	0.67
1:B:69:LEU:HD21	3:B:603:EDO:H12	1.77	0.66
1:B:277:GLN:NE2	1:B:482:ARG:HH12	1.93	0.66
1:A:228:LEU:CD2	5:A:603:GOL:H12	2.27	0.64
1:A:68:ARG:HG2	1:B:329:ILE:HG21	1.79	0.64
1:A:319:THR:HG21	1:A:329:ILE:HD11	1.81	0.63
1:A:139[B]:MSE:SE	1:A:140:LEU:HD23	2.48	0.63
1:A:24:ARG:CG	5:A:609:GOL:H11	2.30	0.61
1:A:1:MSE:HB3	1:A:195:GLU:OE2	2.00	0.61
1:A:24:ARG:HD3	5:A:609:GOL:C1	2.30	0.61
1:B:259:LEU:HD21	1:B:467[B]:LEU:HD21	1.82	0.60
1:B:11:ILE:HD13	1:B:221:LEU:HD23	1.85	0.59
1:B:60:LEU:HD22	1:B:64:TRP:CE2	2.37	0.58
1:A:24:ARG:CD	5:A:609:GOL:H11	2.31	0.58
1:A:116:ARG:HG2	1:A:126:LEU:HD22	1.87	0.57
1:A:344:TYR:O	1:A:347[A]:ARG:HG2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:319:THR:HG21	1:B:329:ILE:HD11	1.86	0.56
1:A:309:ALA:HA	1:A:411:ILE:HD13	1.87	0.56
1:A:355:ARG:NE	6:A:841:HOH:O	2.40	0.54
1:A:25:GLU:OE1	1:B:430:ARG:NH2	2.38	0.53
1:A:329:ILE:HG21	1:B:68:ARG:HG2	1.91	0.53
1:A:260:LEU:HD11	1:A:451:VAL:HB	1.92	0.51
1:A:29:GLU:OE2	1:B:430:ARG:HG2	2.10	0.50
2:A:601:ACT:H3	6:B:764:HOH:O	2.10	0.50
1:A:24:ARG:HD3	5:A:609:GOL:H31	1.94	0.50
1:B:398:SER:O	2:B:602:ACT:CH3	2.60	0.50
1:A:344:TYR:O	1:A:347[B]:ARG:HG2	2.10	0.50
1:B:324:PRO:HA	1:B:327:TYR:CE2	2.47	0.49
1:B:24:ARG:HD3	6:B:711:HOH:O	2.12	0.49
1:A:196:ARG:NH1	6:A:895:HOH:O	2.41	0.49
1:B:368:ARG:HA	1:B:372:GLU:HB2	1.93	0.49
1:A:338[B]:GLU:OE1	1:B:336:THR:OG1	2.22	0.49
1:A:24:ARG:HD3	5:A:609:GOL:C3	2.43	0.49
1:B:259:LEU:CG	1:B:467[B]:LEU:HD21	2.43	0.49
1:A:187:TYR:O	1:A:191:VAL:HG22	2.12	0.49
1:B:259:LEU:CD2	1:B:467[B]:LEU:HD21	2.44	0.48
1:B:408:TRP:HB3	1:B:467[A]:LEU:HD23	1.95	0.48
1:A:25:GLU:HG2	1:B:428:TYR:CZ	2.49	0.48
1:B:398:SER:O	2:B:602:ACT:H3	2.14	0.47
1:A:21:ILE:HG12	5:A:609:GOL:H2	1.95	0.47
1:B:7:VAL:HG22	1:B:9:HIS:CE1	2.50	0.46
1:B:193:TYR:O	1:B:197:VAL:HG23	2.15	0.46
1:B:231:LEU:O	1:B:235:LEU:HG	2.15	0.46
1:A:139[A]:MSE:SE	1:A:345:LEU:HD12	2.65	0.46
1:A:71[B]:GLU:HG2	1:A:73:GLN:NE2	2.31	0.45
1:A:428:TYR:CZ	1:B:25:GLU:HB3	2.52	0.45
1:B:69:LEU:HD11	3:B:603:EDO:H11	1.97	0.45
1:A:117:VAL:O	1:A:189:ARG:NH2	2.49	0.45
1:A:338[B]:GLU:OE1	1:B:338:GLU:HB2	2.16	0.45
1:A:271:LYS:O	1:A:275:LYS:HG3	2.17	0.45
1:B:118:GLU:HB3	1:B:120:HIS:CE1	2.52	0.44
1:A:274:TYR:CZ	1:A:278:LEU:HD22	2.53	0.44
1:A:398[B]:SER:OG	1:A:462:ILE:HG23	2.17	0.44
1:A:14:ASP:OD2	6:A:764:HOH:O	2.21	0.44
1:B:395:GLU:O	1:B:399:GLU:HG3	2.18	0.43
1:B:11:ILE:HD11	6:B:957:HOH:O	2.18	0.43
1:B:75:PHE:HB2	1:B:106:GLY:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:ARG:HG3	1:A:477:ARG:HD3	2.00	0.43
1:B:117:VAL:HG13	1:B:189:ARG:HG3	2.00	0.43
1:A:92:TRP:CH2	1:A:139[B]:MSE:CE	3.02	0.43
1:A:228:LEU:HD21	5:A:603:GOL:H31	2.00	0.42
1:B:405:VAL:HG11	1:B:407:PHE:CZ	2.55	0.42
1:A:347[B]:ARG:HD3	1:A:348:PHE:CZ	2.55	0.42
1:A:75:PHE:HB2	1:A:106:GLY:O	2.20	0.42
1:A:11:ILE:HD12	1:A:218:TYR:HA	2.01	0.41
1:A:225:ARG:HG2	5:A:602:GOL:C3	2.43	0.41
1:B:23:LEU:HD21	1:B:224:GLY:HA3	2.02	0.41
1:A:24:ARG:HG2	5:A:609:GOL:H11	2.01	0.41
1:A:347[B]:ARG:HG3	1:A:348:PHE:CD2	2.56	0.41
1:A:447:TYR:O	1:A:451:VAL:HG23	2.21	0.40
1:A:32:PHE:O	1:A:36:GLN:HG2	2.21	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	467/500 (93%)	459 (98%)	8 (2%)	0	100	100
1	B	454/500 (91%)	440 (97%)	14 (3%)	0	100	100
All	All	921/1000 (92%)	899 (98%)	22 (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	402/418 (96%)	399 (99%)	3 (1%)	87	90
1	B	390/418 (93%)	378 (97%)	12 (3%)	45	44
All	All	792/836 (95%)	777 (98%)	15 (2%)	64	66

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	SER
1	A	457[A]	ARG
1	A	457[B]	ARG
1	B	4	SER
1	B	24	ARG
1	B	60	LEU
1	B	71	GLU
1	B	110	SER
1	B	188	ASP
1	B	292	LYS
1	B	311	LYS
1	B	358	PHE
1	B	363	ARG
1	B	384	LEU
1	B	442	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	306	ASN
1	B	31	GLN
1	B	120	HIS
1	B	277	GLN

5.3.3 RNA ⓘ

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](#)

13 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	ACT	A	601	-	1,3,3	0.81	0	0,3,3	0.00	-
5	GOL	A	602	-	5,5,5	0.87	0	5,5,5	1.45	1 (20%)
5	GOL	A	603	-	5,5,5	0.43	0	5,5,5	0.71	0
3	EDO	A	604	-	3,3,3	0.69	0	2,2,2	0.21	0
2	ACT	A	606	-	1,3,3	2.42	1 (100%)	0,3,3	0.00	-
2	ACT	A	607	-	1,3,3	1.65	0	0,3,3	0.00	-
2	ACT	A	608	-	1,3,3	2.29	1 (100%)	0,3,3	0.00	-
5	GOL	A	609	-	5,5,5	0.67	0	5,5,5	1.17	0
5	GOL	A	610	-	5,5,5	0.31	0	5,5,5	0.81	0
2	ACT	B	602	-	1,3,3	1.26	0	0,3,3	0.00	-
3	EDO	B	603	-	3,3,3	0.50	0	2,2,2	0.20	0
4	MLI	B	605	-	0,6,6	0.00	-	0,7,7	0.00	-
3	EDO	B	606	-	3,3,3	0.44	0	2,2,2	0.47	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
2	ACT	A	601	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GOL	A	602	-	-	0/4/4/4	0/0/0/0
5	GOL	A	603	-	-	0/4/4/4	0/0/0/0
3	EDO	A	604	-	-	0/1/1/1	0/0/0/0
2	ACT	A	606	-	-	0/0/0/0	0/0/0/0
2	ACT	A	607	-	-	0/0/0/0	0/0/0/0
2	ACT	A	608	-	-	0/0/0/0	0/0/0/0
5	GOL	A	609	-	-	0/4/4/4	0/0/0/0
5	GOL	A	610	-	-	0/4/4/4	0/0/0/0
2	ACT	B	602	-	-	0/0/0/0	0/0/0/0
3	EDO	B	603	-	-	0/1/1/1	0/0/0/0
4	MLI	B	605	-	-	0/0/4/4	0/0/0/0
3	EDO	B	606	-	-	0/1/1/1	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	608	ACT	CH3-C	2.29	1.51	1.48
2	A	606	ACT	CH3-C	2.42	1.51	1.48

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	602	GOL	C3-C2-C1	2.47	121.35	111.52

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	ACT	1	0
5	A	602	GOL	2	0
5	A	603	GOL	2	0
5	A	609	GOL	8	0
2	B	602	ACT	2	0
3	B	603	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	460/500 (92%)	0.54	36 (7%)	14 14	16, 29, 50, 59	0
1	B	452/500 (90%)	0.67	40 (8%)	11 11	13, 29, 55, 86	1 (0%)
All	All	912/1000 (91%)	0.61	76 (8%)	12 12	13, 29, 53, 86	1 (0%)

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	427	PRO	10.0
1	A	176	LEU	8.8
1	B	2	PHE	7.7
1	A	255	ASP	7.3
1	B	7	VAL	6.7
1	A	428	TYR	6.3
1	B	238	LYS	6.2
1	A	182	LYS	5.4
1	A	431	TYR	5.4
1	B	5	ALA	5.3
1	B	124	ALA	5.2
1	B	157	LYS	5.1
1	A	177	SER	5.0
1	A	426	THR	4.8
1	B	4	SER	4.4
1	A	436	GLU	4.4
1	B	496	LYS	4.4
1	A	435	GLU	4.2
1	A	439	ARG	4.1
1	A	432	LYS	4.0
1	B	379	ALA	3.8
1	B	237	THR	3.8
1	B	236	ALA	3.8
1	A	433	ILE	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	434	THR	3.7
1	B	161	GLU	3.7
1	A	76	LEU	3.7
1	B	121	ILE	3.6
1	B	196	ARG	3.5
1	B	3	GLU	3.4
1	B	192	HIS	3.2
1	B	8	GLY	3.1
1	B	117	VAL	3.1
1	A	438	TRP	2.9
1	A	496	LYS	2.9
1	A	185	GLU	2.9
1	B	119	GLY	2.8
1	B	128	GLN	2.8
1	A	257	ARG	2.8
1	B	260	LEU	2.8
1	A	180	ASP	2.7
1	B	439	ARG	2.7
1	A	179	LEU	2.7
1	A	329	ILE	2.6
1	B	188	ASP	2.6
1	A	259	LEU	2.6
1	B	163	LEU	2.6
1	B	162	ARG	2.5
1	A	256	ASN	2.5
1	B	118	GLU	2.5
1	B	329	ILE	2.5
1	A	442	ASP	2.4
1	A	114	TYR	2.4
1	A	161	GLU	2.4
1	B	315	ILE	2.4
1	A	425	LYS	2.3
1	B	160	LYS	2.3
1	B	193	TYR	2.3
1	B	189	ARG	2.3
1	B	114	TYR	2.2
1	B	348	PHE	2.2
1	A	260	LEU	2.2
1	A	5	ALA	2.2
1	B	360	ILE	2.1
1	B	186	VAL	2.1
1	A	181	TRP	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	18	LYS	2.1
1	B	123	GLU	2.1
1	A	184	SER	2.1
1	A	437	ASP	2.0
1	B	442	ASP	2.0
1	A	183	GLN	2.0
1	A	124	ALA	2.0
1	A	8	GLY	2.0
1	B	310	GLY	2.0
1	B	435	GLU	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
5	GOL	A	609	6/6	0.88	0.48	13.81	42,52,53,54	0
5	GOL	A	602	6/6	0.74	0.40	10.25	52,57,58,58	0
5	GOL	A	603	6/6	0.89	0.34	9.43	53,56,57,58	0
2	ACT	A	608	4/4	0.86	0.27	8.10	56,56,57,57	0
2	ACT	B	602	4/4	0.90	0.28	6.25	56,56,56,57	0
3	EDO	B	603	4/4	0.87	0.18	0.71	59,61,62,64	0
5	GOL	A	610	6/6	0.84	0.16	0.38	69,71,71,71	0
3	EDO	B	606	4/4	0.85	0.16	0.27	54,54,55,58	0
4	MLI	B	605	7/7	0.89	0.13	-1.51	37,47,52,60	0
2	ACT	A	601	4/4	0.94	0.13	-2.45	62,62,62,63	0
3	EDO	A	604	4/4	0.89	0.33	-	57,59,59,60	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	ACT	A	606	4/4	0.81	0.16	-	67,67,67,67	0
2	ACT	A	607	4/4	0.89	0.21	-	55,55,56,56	0

6.5 Other polymers [i](#)

There are no such residues in this entry.