



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:30 PM GMT

PDB ID : 4P3Z
Title : Chlamydia pneumoniae CopN (D29 construct)
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Deposited on : 2014-03-10
Resolution : 1.77 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

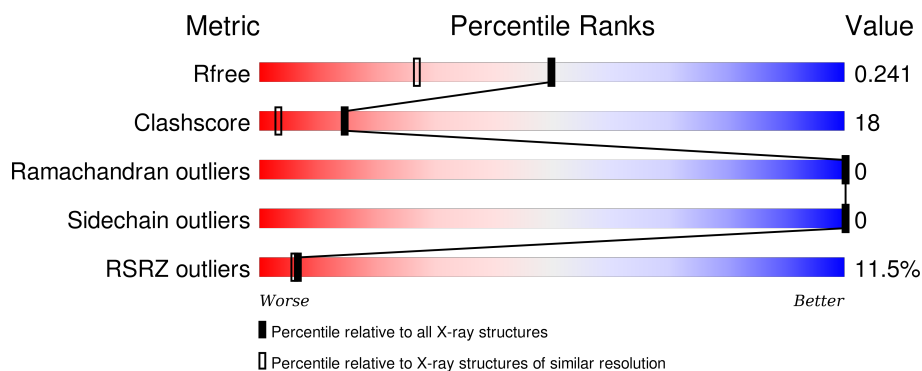
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.77 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	6655 (1.80-1.76)
Clashscore	102246	7658 (1.80-1.76)
Ramachandran outliers	100387	7570 (1.80-1.76)
Sidechain outliers	100360	7569 (1.80-1.76)
RSRZ outliers	91569	6671 (1.80-1.76)

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	425	

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	SO4	A	402	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	403	-	-	-	X
2	SO4	A	404	-	-	-	X
3	GOL	A	406	-	-	-	X
3	GOL	A	407	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4768 atoms, of which 2287 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 CopN.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	292	Total	C	H	N	O	S	0	4	0
			4564	1433	2271	390	463	7			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-25	MET	-	initiating methionine	UNP Q9Z8L4
A	-24	LYS	-	expression tag	UNP Q9Z8L4
A	-23	HIS	-	expression tag	UNP Q9Z8L4
A	-22	HIS	-	expression tag	UNP Q9Z8L4
A	-21	HIS	-	expression tag	UNP Q9Z8L4
A	-20	HIS	-	expression tag	UNP Q9Z8L4
A	-19	HIS	-	expression tag	UNP Q9Z8L4
A	-18	HIS	-	expression tag	UNP Q9Z8L4
A	-17	PRO	-	expression tag	UNP Q9Z8L4
A	-16	MET	-	expression tag	UNP Q9Z8L4
A	-15	SER	-	expression tag	UNP Q9Z8L4
A	-14	ASP	-	expression tag	UNP Q9Z8L4
A	-13	TYR	-	expression tag	UNP Q9Z8L4
A	-12	ASP	-	expression tag	UNP Q9Z8L4
A	-11	ILE	-	expression tag	UNP Q9Z8L4
A	-10	PRO	-	expression tag	UNP Q9Z8L4
A	-9	THR	-	expression tag	UNP Q9Z8L4
A	-8	THR	-	expression tag	UNP Q9Z8L4
A	-7	GLU	-	expression tag	UNP Q9Z8L4
A	-6	ASN	-	expression tag	UNP Q9Z8L4
A	-5	LEU	-	expression tag	UNP Q9Z8L4
A	-4	TYR	-	expression tag	UNP Q9Z8L4
A	-3	PHE	-	expression tag	UNP Q9Z8L4
A	-2	GLN	-	expression tag	UNP Q9Z8L4
A	-1	GLY	-	expression tag	UNP Q9Z8L4
A	0	ALA	-	expression tag	UNP Q9Z8L4

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).

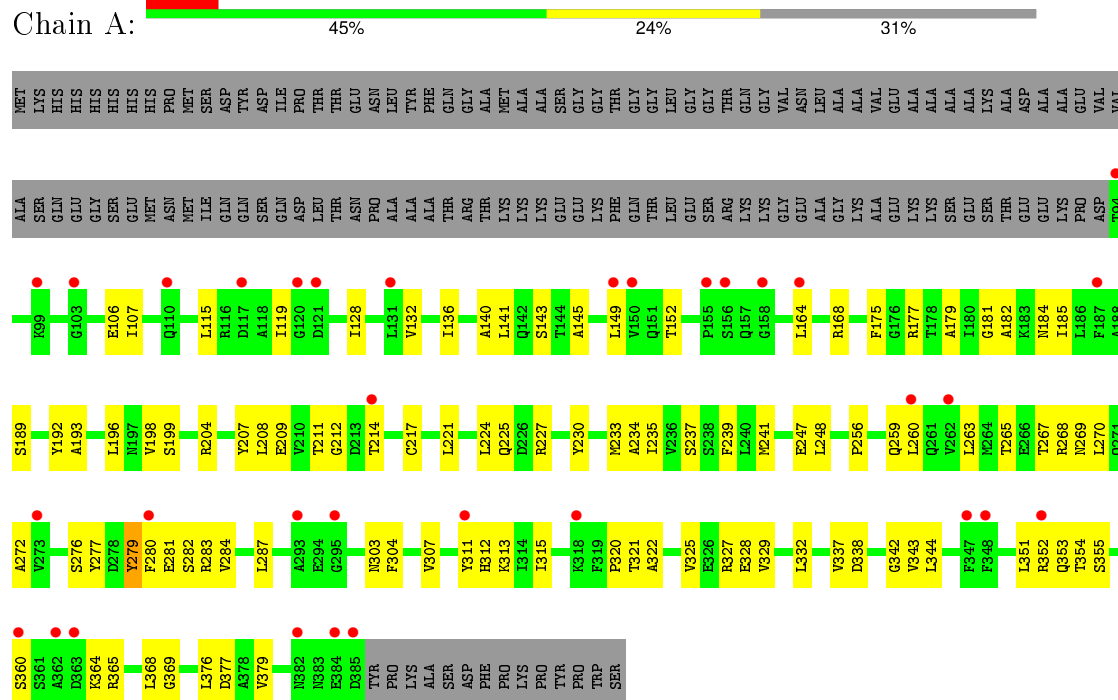


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	151	Total	O	0	0
			151	151		

- Molecule 1: CopN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	62.70 Å 62.70 Å 85.70 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	16.29 – 1.77 16.29 – 1.77	Depositor EDS
% Data completeness (in resolution range)	99.6 (16.29-1.77) 99.6 (16.29-1.77)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 1.77 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.180 , 0.235 0.188 , 0.241	Depositor DCC
R_{free} test set	1668 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	28.8	Xtriage
Anisotropy	0.067	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.49 , 34.0	EDS
Estimated twinning fraction	0.021 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 33432 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4768	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.00% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MLZ, GOL, 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	0.58	1/2287 (0.0%)	0.52	0/3095

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	279	TYR	CE1-CZ	-5.56	1.31	1.38

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2293	2271	2280	83	1
2	A	25	0	0	4	0
3	A	12	16	16	2	0
4	A	151	0	0	1	0
All	All	2481	2287	2296	83	1

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 18.

All (83) 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:221:LEU:O	1:A:225[A]:GLN:HG3	1.83	0.78
1:A:272:ALA:HB1	1:A:355:SER:HB3	1.71	0.72
1:A:268:ARG:NH1	2:A:402:SO4:O4	2.23	0.70
1:A:224:LEU:HD12	1:A:270:LEU:HD21	1.75	0.67
1:A:248:LEU:HD21	1:A:260:LEU:HB2	1.79	0.65
1:A:337:VAL:HG23	1:A:379:VAL:HG12	1.78	0.65
1:A:282:SER:OG	1:A:283:ARG:NH1	2.31	0.64
1:A:311:TYR:OH	1:A:351:LEU:HA	1.98	0.63
1:A:311:TYR:OH	1:A:354:THR:OG1	1.92	0.62
1:A:196:LEU:HD21	1:A:235:ILE:HB	1.82	0.61
1:A:211:THR:HA	1:A:263:LEU:HD13	1.83	0.60
1:A:106:GLU:OE1	1:A:204:ARG:NH1	2.34	0.60
1:A:284:VAL:HG21	1:A:304:PHE:HA	1.81	0.60
1:A:182:ALA:HB2	1:A:208:LEU:HD11	1.81	0.60
1:A:307:VAL:HA	1:A:332:LEU:HD21	1.85	0.58
1:A:107:ILE:HG23	1:A:141:LEU:HD22	1.87	0.57
1:A:128:ILE:HG21	1:A:164:LEU:HD23	1.87	0.57
1:A:321:THR:O	1:A:325:VAL:HG23	2.06	0.56
1:A:221:LEU:HG	1:A:225[A]:GLN:NE2	2.21	0.55
1:A:352:ARG:NH2	2:A:404:SO4:O1	2.40	0.54
1:A:185:ILE:HD13	1:A:207:TYR:CG	2.43	0.54
1:A:140:ALA:HB2	1:A:182:ALA:CB	2.37	0.54
1:A:119:ILE:CD1	1:A:149:LEU:HD13	2.38	0.53
1:A:128:ILE:HD13	1:A:164:LEU:HA	1.92	0.52
1:A:214:THR:HG22	4:A:639:HOH:O	2.10	0.51
1:A:230:TYR:HA	1:A:233:MET:HE3	1.93	0.50
1:A:192:TYR:CG	1:A:239:PHE:HB2	2.47	0.50
1:A:177:ARG:O	1:A:181:GLY:N	2.42	0.50
1:A:193:ALA:HB1	1:A:198:VAL:O	2.12	0.50
1:A:140:ALA:HB2	1:A:182:ALA:HB3	1.93	0.49
1:A:217:CYS:HB2	1:A:269:ASN:HB2	1.94	0.49
1:A:281:GLU:HA	1:A:304:PHE:CG	2.47	0.49
1:A:281:GLU:HA	1:A:304:PHE:CB	2.43	0.48
1:A:351:LEU:HD21	1:A:365:ARG:O	2.13	0.48
1:A:143:SER:HB2	1:A:179:ALA:HB1	1.94	0.48
1:A:276:SER:OG	1:A:312:HIS:NE2	2.36	0.48
1:A:354:THR:O	1:A:365:ARG:NH1	2.41	0.48
1:A:344:LEU:HB3	1:A:376:LEU:HG	1.96	0.48
1:A:175:PHE:CB	1:A:179:ALA:HB2	2.44	0.47
1:A:211:THR:HA	1:A:263:LEU:CD1	2.45	0.47
1:A:313:MLZ:HCM2	1:A:328:GLU:OE2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:LEU:HD21	1:A:132:VAL:HG22	1.96	0.47
1:A:185:ILE:HG21	1:A:207:TYR:CZ	2.50	0.47
1:A:207:TYR:CZ	1:A:211:THR:HG21	2.49	0.47
1:A:276:SER:CB	1:A:312:HIS:NE2	2.78	0.47
1:A:115:LEU:HD21	1:A:132:VAL:HA	1.97	0.47
1:A:322:ALA:H	3:A:406:GOL:H32	1.80	0.47
1:A:287:LEU:HD11	1:A:342:GLY:HA3	1.98	0.46
1:A:149:LEU:HB3	1:A:164:LEU:HD22	1.99	0.45
1:A:351:LEU:O	1:A:365:ARG:NH1	2.50	0.45
1:A:177:ARG:NH2	1:A:247:GLU:OE1	2.32	0.45
1:A:189:SER:HA	1:A:239:PHE:CE1	2.51	0.45
1:A:208:LEU:O	1:A:212:GLY:N	2.50	0.45
1:A:265:THR:HG23	2:A:402:SO4:O2	2.18	0.44
1:A:329:VAL:HG21	1:A:344:LEU:CD2	2.48	0.43
1:A:315:ILE:HG22	1:A:368:LEU:HD13	1.99	0.43
1:A:241:MET:SD	1:A:267:THR:HB	2.58	0.43
1:A:217:CYS:SG	1:A:270:LEU:HA	2.58	0.43
1:A:344:LEU:HD12	1:A:379:VAL:HG21	2.01	0.43
1:A:136:ILE:CD1	1:A:145:ALA:HB2	2.48	0.43
1:A:136:ILE:CG2	1:A:141:LEU:HB2	2.49	0.43
1:A:199:SER:N	2:A:405:SO4:O1	2.48	0.42
1:A:234:ALA:O	1:A:237[A]:SER:OG	2.31	0.42
1:A:185:ILE:HD11	1:A:204:ARG:HA	2.01	0.42
1:A:230:TYR:CE1	1:A:277:TYR:CD2	3.07	0.42
1:A:337:VAL:HG13	1:A:338:ASP:N	2.34	0.42
1:A:332:LEU:HD23	1:A:343:VAL:HG11	2.02	0.42
1:A:276:SER:O	1:A:279:TYR:HB3	2.19	0.42
1:A:209:GLU:CD	1:A:227:ARG:HH22	2.23	0.42
1:A:351:LEU:CD2	1:A:369:GLY:CA	2.97	0.42
1:A:107:ILE:CG2	1:A:141:LEU:HD22	2.49	0.42
1:A:303:ASN:O	1:A:307:VAL:HG23	2.20	0.42
1:A:184:ASN:HB3	1:A:247:GLU:HG2	2.02	0.41
1:A:280:PHE:O	1:A:284:VAL:HG23	2.20	0.41
1:A:360:SER:OG	1:A:364:MLZ:HG2	2.20	0.41
1:A:149:LEU:HA	1:A:152:THR:HG22	2.02	0.41
1:A:320:PRO:HG3	1:A:368:LEU:HD21	2.02	0.41
1:A:327:ARG:HA	1:A:327:ARG:NE	2.36	0.41
1:A:312:HIS:HA	1:A:315:ILE:HG12	2.02	0.41
1:A:351:LEU:CD2	1:A:369:GLY:N	2.84	0.41
1:A:279:TYR:CE1	1:A:353:GLN:HG3	2.57	0.40
1:A:256:PRO:HB2	1:A:259:GLN:HB3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:ALA:H	3:A:406:GOL:C3	2.33	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:ARG:HH12	1:A:377:ASP:OD1[2_345]	1.58	0.02

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	290/425 (68%)	287 (99%)	3 (1%)	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	251/350 (72%)	251 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	MLZ	A	249	1	8,9,10	0.84	0	7,9,11	1.91	1 (14%)
1	MLZ	A	313	1	8,9,10	0.73	0	7,9,11	1.61	1 (14%)
1	MLZ	A	324	1	8,9,10	0.75	0	7,9,11	1.63	2 (28%)
1	MLZ	A	364	1	8,9,10	0.69	0	7,9,11	1.44	2 (28%)

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
1	MLZ	A	249	1	-	0/6/8/10	0/0/0/0
1	MLZ	A	313	1	-	0/6/8/10	0/0/0/0
1	MLZ	A	324	1	-	0/6/8/10	0/0/0/0
1	MLZ	A	364	1	-	0/6/8/10	0/0/0/0

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	364	MLZ	O-C-CA	-2.40	119.23	125.49
1	A	324	MLZ	O-C-CA	-2.20	119.77	125.49
1	A	364	MLZ	CM-NZ-CE	2.66	120.00	112.23
1	A	324	MLZ	CM-NZ-CE	3.21	121.62	112.23
1	A	313	MLZ	CM-NZ-CE	3.37	122.10	112.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	249	MLZ	CM-NZ-CE	3.87	123.56	112.23

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
1	A	313	MLZ	1	0
1	A	364	MLZ	1	0

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

7 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	SO4	A	401	-	4,4,4	0.25	0	6,6,6	0.07	0
2	SO4	A	402	-	4,4,4	0.28	0	6,6,6	0.07	0
2	SO4	A	403	-	4,4,4	0.25	0	6,6,6	0.07	0
2	SO4	A	404	-	4,4,4	0.23	0	6,6,6	0.06	0
2	SO4	A	405	-	4,4,4	0.18	0	6,6,6	0.08	0
3	GOL	A	406	-	5,5,5	0.15	0	5,5,5	0.66	0
3	GOL	A	407	-	5,5,5	0.19	0	5,5,5	0.19	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	SO4	A	401	-	-	0/0/0/0	0/0/0/0
2	SO4	A	402	-	-	0/0/0/0	0/0/0/0
2	SO4	A	403	-	-	0/0/0/0	0/0/0/0
2	SO4	A	404	-	-	0/0/0/0	0/0/0/0
2	SO4	A	405	-	-	0/0/0/0	0/0/0/0
3	GOL	A	406	-	-	0/4/4/4	0/0/0/0
3	GOL	A	407	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	402	SO4	2	0
2	A	404	SO4	1	0
2	A	405	SO4	1	0
3	A	406	GOL	2	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	288/425 (67%)	0.89	33 (11%) 6 6	22, 30, 54, 77	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	103	GLY	7.1
1	A	385	ASP	5.9
1	A	94	THR	5.6
1	A	384	GLU	5.1
1	A	156	SER	4.9
1	A	155	PRO	4.1
1	A	311	TYR	3.8
1	A	362	ALA	3.3
1	A	110	GLN	3.2
1	A	363	ASP	3.2
1	A	131	LEU	3.1
1	A	149	LEU	3.1
1	A	117	ASP	2.7
1	A	121	ASP	2.7
1	A	187[A]	PHE	2.6
1	A	295	GLY	2.6
1	A	360	SER	2.6
1	A	318	LYS	2.6
1	A	347	PHE	2.5
1	A	150	VAL	2.5
1	A	164	LEU	2.5
1	A	352	ARG	2.3
1	A	293	ALA	2.3
1	A	382	ASN	2.2
1	A	120	GLY	2.2
1	A	260	LEU	2.2
1	A	348	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	99	LYS	2.2
1	A	214	THR	2.2
1	A	273	VAL	2.0
1	A	158	GLY	2.0
1	A	280	PHE	2.0
1	A	262	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	MLZ	A	249	10/11	0.84	0.21	-	24,37,50,50	0
1	MLZ	A	364	10/11	0.92	0.15	-	35,55,66,66	0
1	MLZ	A	313	10/11	0.90	0.18	-	26,51,65,65	0
1	MLZ	A	324	10/11	0.92	0.11	-	24,37,45,45	0

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
3	GOL	A	406	6/6	0.57	0.39	7.31	46,60,74,75	0
2	SO4	A	403	5/5	0.90	0.18	3.56	60,62,63,63	5
2	SO4	A	404	5/5	0.78	0.30	2.97	63,64,65,66	5
3	GOL	A	407	6/6	0.53	0.41	2.23	76,92,93,94	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	SO4	A	401	5/5	0.97	0.21	1.16	61,62,63,63	5
2	SO4	A	402	5/5	0.97	0.20	-	50,53,57,58	5
2	SO4	A	405	5/5	0.89	0.33	-	66,68,68,69	0

6.5 Other polymers [i](#)

There are no such residues in this entry.