



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 19, 2022 – 08:21 AM JST

PDB ID : 7FF0
Title : Pseudomonas aeruginosa Virulence Factor Regulator with cAMP ligand and auranofin
Authors : Chew, B.L.A.; Luo, D.
Deposited on : 2021-07-22
Resolution : 2.60 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	:	4.02b-467
Mogul	:	?? (??), CSD ??CSD?? (????)
Xtriage (Phenix)	:	1.13
EDS	:	2.29
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.29

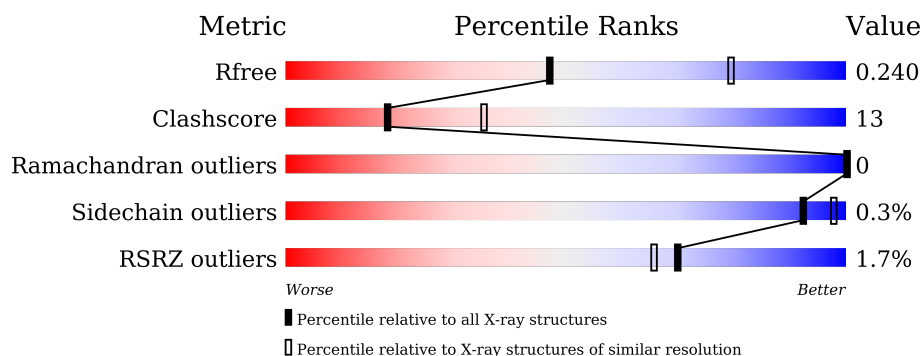
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	213	<div> <div>2%</div> <div> <div></div> <div>71%</div> <div>25%</div> <div>..</div> </div> </div>
1	B	213	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>15%</div> <div>6%</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
3	SO4	A	303	-	-	X	-
3	SO4	A	304	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6456 atoms, of which 3099 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.

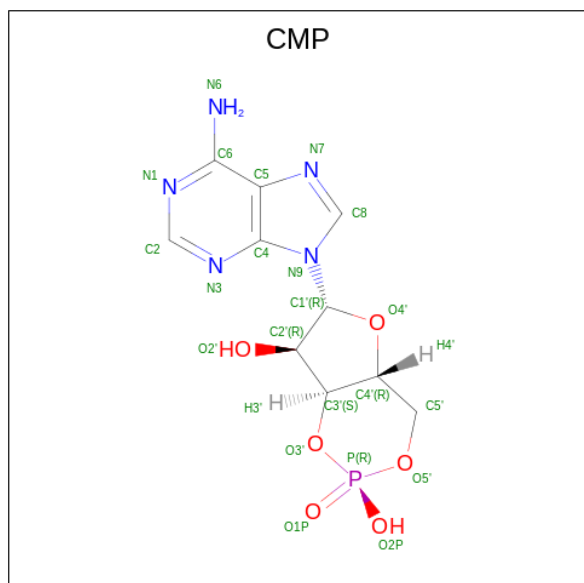
- Molecule 1 is a protein called cAMP-activated global transcriptional regulator Vfr.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	206	Total	C	H	N	O	S	0	0	0
			3174	997	1581	282	303	11			
1	B	200	Total	C	H	N	O	S	0	0	0
			3024	957	1496	274	286	11			

There are 6 discrepancies between the modelled and reference sequences:

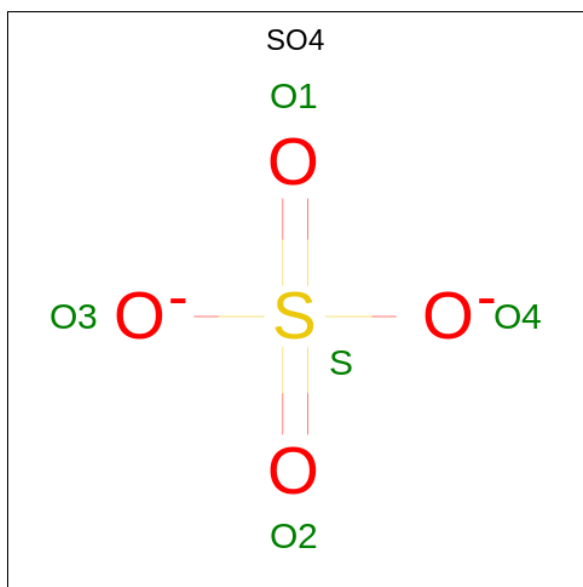
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	THR	-	expression tag	UNP P55222
A	3	GLY	-	expression tag	UNP P55222
A	4	GLY	-	expression tag	UNP P55222
B	2	THR	-	expression tag	UNP P55222
B	3	GLY	-	expression tag	UNP P55222
B	4	GLY	-	expression tag	UNP P55222

- Molecule 2 is ADENOSINE-3',5'-CYCLIC-MONOPHOSPHATE (three-letter code: CMP) (formula: C₁₀H₁₂N₅O₆P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	P	
			33	10	11	5	6	1	
2	B	1	Total	C	H	N	O	P	
			33	10	11	5	6	1	

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 4 is GOLD ION (three-letter code: AU) (formula: Au) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Au		
			1	1	0	0

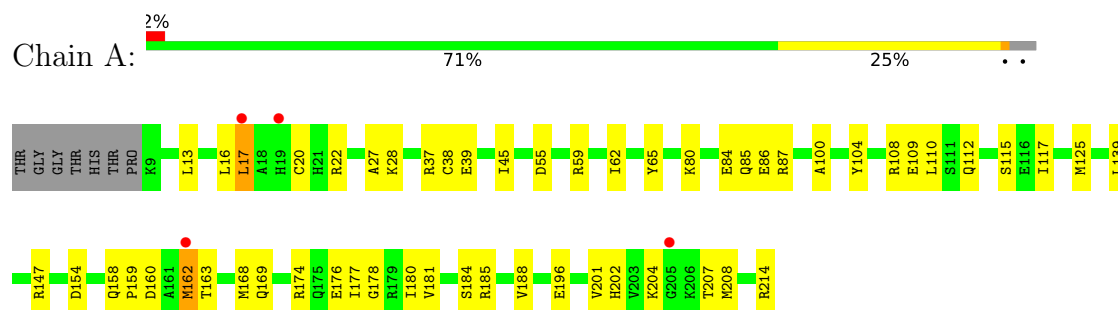
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	100	Total 100	O 100	0	0
5	B	66	Total 66	O 66	0	0

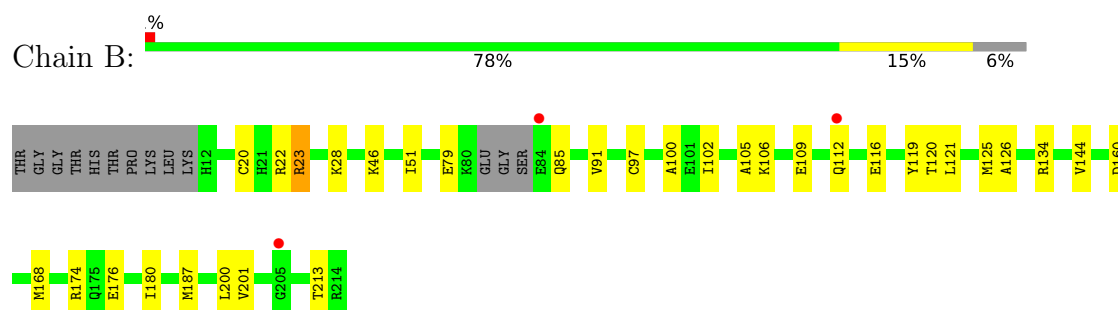
3 Residue-property plots [i](#)

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: cAMP-activated global transcriptional regulator Vfr



- Molecule 1: cAMP-activated global transcriptional regulator Vfr



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	135.60Å 49.64Å 63.91Å 90.00° 100.37° 90.00°	Depositor
Resolution (Å)	42.12 – 2.60 42.12 – 2.60	Depositor EDS
% Data completeness (in resolution range)	93.0 (42.12-2.60) 90.1 (42.12-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.207 , 0.239 0.207 , 0.240	Depositor DCC
R_{free} test set	609 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	0.277	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 62.3	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	0.91	EDS
Total number of atoms	6456	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.69% 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: AU, CMP, 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	0.59	3/1612 (0.2%)	0.86	5/2166 (0.2%)
1	B	0.42	0/1547	0.78	1/2080 (0.0%)
All	All	0.52	3/3159 (0.1%)	0.83	6/4246 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	162	MET	CB-CG	11.98	1.89	1.51
1	A	162	MET	CG-SD	8.71	2.03	1.81
1	A	162	MET	CA-CB	5.47	1.66	1.53

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	162	MET	CA-CB-CG	-10.47	95.50	113.30
1	B	23	ARG	CG-CD-NE	-9.54	91.77	111.80
1	A	17	LEU	CA-CB-CG	-8.26	96.30	115.30
1	A	162	MET	N-CA-CB	5.80	121.03	110.60
1	A	110	LEU	CA-CB-CG	-5.50	102.65	115.30
1	A	208	MET	CG-SD-CE	5.26	108.61	100.20

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	1593	1581	1581	56	0
1	B	1528	1496	1496	26	2
2	A	22	11	11	1	0
2	B	22	11	11	1	0
3	A	15	0	0	7	2
3	B	10	0	0	0	0
4	A	1	0	0	0	0
5	A	100	0	0	9	2
5	B	66	0	0	7	2
All	All	3357	3099	3099	84	4

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 13.

All (84) 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)
2:A:301:CMP:H2	2:A:301:CMP:C2	0.97	1.50
2:B:301:CMP:H2	2:B:301:CMP:C2	0.97	1.50
1:A:162:MET:CB	1:A:162:MET:CG	1.89	1.45
1:A:162:MET:CG	1:A:162:MET:SD	2.03	1.45
1:A:37:ARG:HD3	1:A:39:GLU:HG3	1.09	1.08
1:A:37:ARG:HD3	1:A:39:GLU:CG	1.88	1.04
1:A:37:ARG:CD	1:A:39:GLU:HG3	1.93	0.99
1:A:162:MET:SD	5:A:406:HOH:O	2.21	0.96
1:A:176:GLU:OE1	5:A:401:HOH:O	1.89	0.90
1:B:116:GLU:O	5:B:401:HOH:O	1.91	0.89
3:A:303:SO4:O2	5:A:402:HOH:O	1.89	0.89
1:B:28:LYS:HE3	1:B:28:LYS:HA	1.57	0.87
3:A:303:SO4:S	5:A:402:HOH:O	2.34	0.83
1:A:115:SER:OG	1:B:112:GLN:O	1.98	0.80
1:A:174:ARG:NH2	3:A:304:SO4:O3	2.18	0.77
1:A:162:MET:CG	1:A:162:MET:CA	2.63	0.75
1:A:16:LEU:HD12	1:A:117:ILE:HG21	1.67	0.75
1:B:134:ARG:NH1	5:B:402:HOH:O	2.08	0.74
1:B:174:ARG:NH1	5:B:403:HOH:O	2.20	0.73
1:B:79:GLU:HG3	1:B:85:GLN:HB2	1.72	0.70
1:A:185:ARG:HA	1:A:188:VAL:HG22	1.72	0.69
1:A:37:ARG:NH2	3:A:303:SO4:O3	2.25	0.69
1:A:84:GLU:OE1	5:A:403:HOH:O	2.13	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:162:MET:CG	1:A:162:MET:CE	2.74	0.65
1:B:120:THR:N	5:B:401:HOH:O	2.01	0.64
1:A:204:LYS:HB2	1:A:207:THR:O	1.98	0.64
1:A:45:ILE:HD11	1:A:100:ALA:HB2	1.79	0.62
1:A:16:LEU:CD1	1:A:117:ILE:HG21	2.29	0.61
1:A:80:LYS:O	5:A:404:HOH:O	2.16	0.60
1:B:105:ALA:O	1:B:109:GLU:HG2	2.02	0.59
1:A:37:ARG:HG2	1:A:38:CYS:N	2.19	0.58
1:A:159:PRO:O	1:A:160:ASP:HB2	2.03	0.57
1:A:185:ARG:HA	1:A:188:VAL:CG2	2.36	0.56
1:A:201:VAL:HG22	1:A:202:HIS:N	2.22	0.55
1:A:65:TYR:OH	1:A:176:GLU:OE2	2.21	0.55
1:B:28:LYS:HA	1:B:28:LYS:CE	2.32	0.54
1:A:196:GLU:HG2	1:A:201:VAL:HG13	1.90	0.54
1:A:37:ARG:HD3	1:A:39:GLU:CD	2.27	0.53
1:B:102:ILE:HD11	1:B:106:LYS:CB	2.39	0.53
1:A:163:THR:HG23	1:A:214:ARG:HD2	1.90	0.52
1:B:144:VAL:HG11	1:B:187:MET:SD	2.49	0.52
1:A:178:GLY:HA2	1:A:188:VAL:HG11	1.90	0.52
1:A:37:ARG:NH2	1:A:39:GLU:OE2	2.43	0.52
1:A:104:TYR:O	1:A:108:ARG:HG3	2.10	0.52
1:A:13:LEU:O	1:A:17:LEU:HD12	2.09	0.51
1:B:168:MET:HG2	1:B:213:THR:CG2	2.39	0.51
1:A:20:CYS:HB3	1:A:100:ALA:HB1	1.93	0.50
1:B:116:GLU:C	5:B:401:HOH:O	2.43	0.50
1:B:176:GLU:O	1:B:180:ILE:HG13	2.12	0.49
1:B:121:LEU:O	1:B:125:MET:HG3	2.13	0.48
1:A:162:MET:HG3	1:A:169:GLN:O	2.14	0.48
1:B:20:CYS:O	1:B:22:ARG:NH1	2.46	0.48
1:B:200:LEU:HG	1:B:201:VAL:HG13	1.95	0.48
1:A:169:GLN:NE2	5:A:406:HOH:O	2.40	0.48
1:A:55:ASP:OD2	1:A:59:ARG:HB3	2.14	0.48
1:B:119:TYR:N	5:B:401:HOH:O	2.47	0.47
1:A:37:ARG:O	3:A:303:SO4:O4	2.32	0.47
1:B:168:MET:HG2	1:B:213:THR:HG21	1.97	0.47
1:B:46:LYS:O	1:B:97:CYS:HA	2.15	0.46
1:A:22:ARG:HD3	1:A:45:ILE:HD13	1.98	0.46
1:A:174:ARG:HH21	1:A:174:ARG:HG3	1.81	0.46
1:A:174:ARG:HG2	3:A:304:SO4:O2	2.17	0.45
1:A:62:ILE:CD1	1:A:176:GLU:HG3	2.46	0.45
1:A:85:GLN:HG3	1:A:86:GLU:N	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:GLU:O	1:A:112:GLN:HB3	2.17	0.44
1:A:125:MET:HE1	1:B:126:ALA:HB2	1.99	0.44
1:A:185:ARG:CA	1:A:188:VAL:HG22	2.46	0.44
1:A:177:ILE:O	1:A:181:VAL:HG22	2.18	0.44
1:A:139:LEU:HA	1:A:147:ARG:HD3	1.98	0.44
1:A:196:GLU:CG	1:A:201:VAL:HG13	2.48	0.43
1:B:51:ILE:HD12	1:B:91:VAL:HG22	1.99	0.43
1:B:160:ASP:OD1	1:B:160:ASP:N	2.49	0.42
1:A:87:ARG:NH1	5:A:402:HOH:O	2.23	0.42
1:A:176:GLU:O	1:A:180:ILE:HG13	2.20	0.42
1:A:162:MET:O	1:A:168:MET:HA	2.19	0.42
1:A:27:ALA:O	1:A:28:LYS:HB2	2.18	0.42
1:A:162:MET:CG	1:A:162:MET:N	2.83	0.42
1:A:174:ARG:N	3:A:304:SO4:O2	2.51	0.42
1:A:154:ASP:O	1:A:158:GLN:HG2	2.19	0.41
1:B:213:THR:O	1:B:213:THR:HG23	2.19	0.41
1:A:22:ARG:NH1	5:A:417:HOH:O	2.54	0.41
1:B:119:TYR:HB3	5:B:401:HOH:O	2.20	0.41
1:A:178:GLY:CA	1:A:188:VAL:HG11	2.52	0.40
1:B:20:CYS:HB3	1:B:100:ALA:HB1	2.03	0.40

All (4) 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:B:23:ARG:HH22	3:A:303:SO4:O3[1_565]	1.02	0.58
1:B:23:ARG:NH2	3:A:303:SO4:O3[1_565]	1.81	0.39
5:A:415:HOH:O	5:B:455:HOH:O[1_554]	1.92	0.28
5:A:495:HOH:O	5:B:448:HOH:O[4_748]	1.96	0.24

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	204/213 (96%)	196 (96%)	8 (4%)	0	100	100
1	B	196/213 (92%)	189 (96%)	7 (4%)	0	100	100
All	All	400/426 (94%)	385 (96%)	15 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	167/182 (92%)	166 (99%)	1 (1%)	86	95
1	B	156/182 (86%)	156 (100%)	0	100	100
All	All	323/364 (89%)	322 (100%)	1 (0%)	92	98

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	184	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	157	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 1 is monoatomic - leaving 7 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

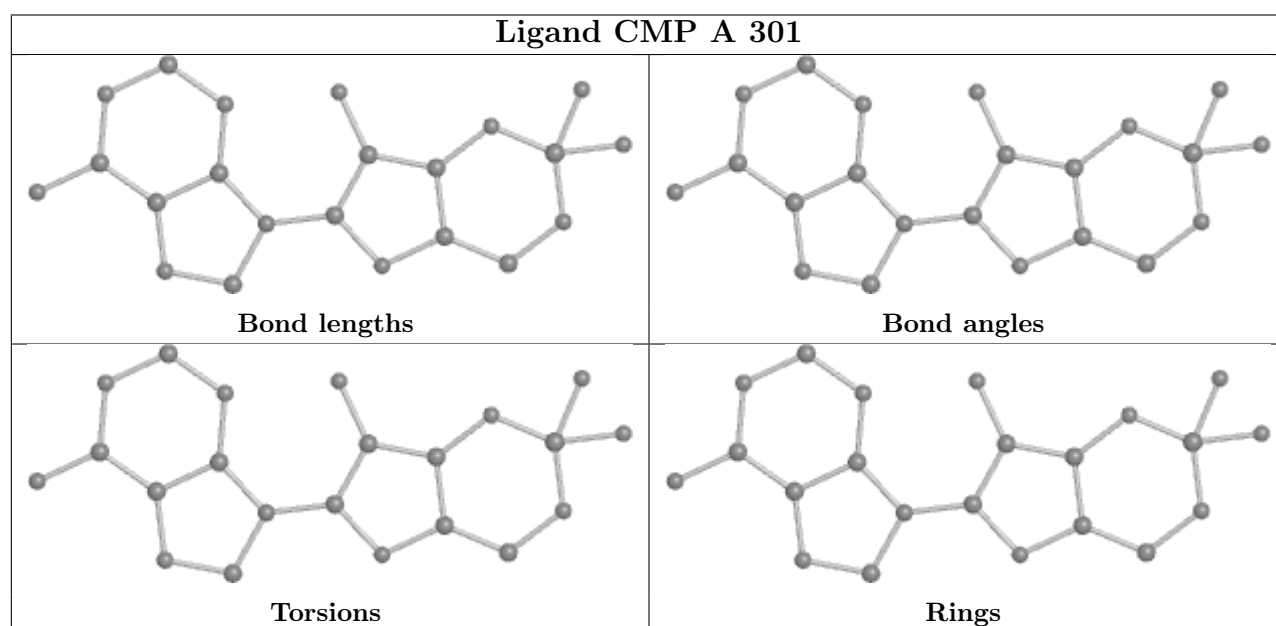
There are no chirality outliers.

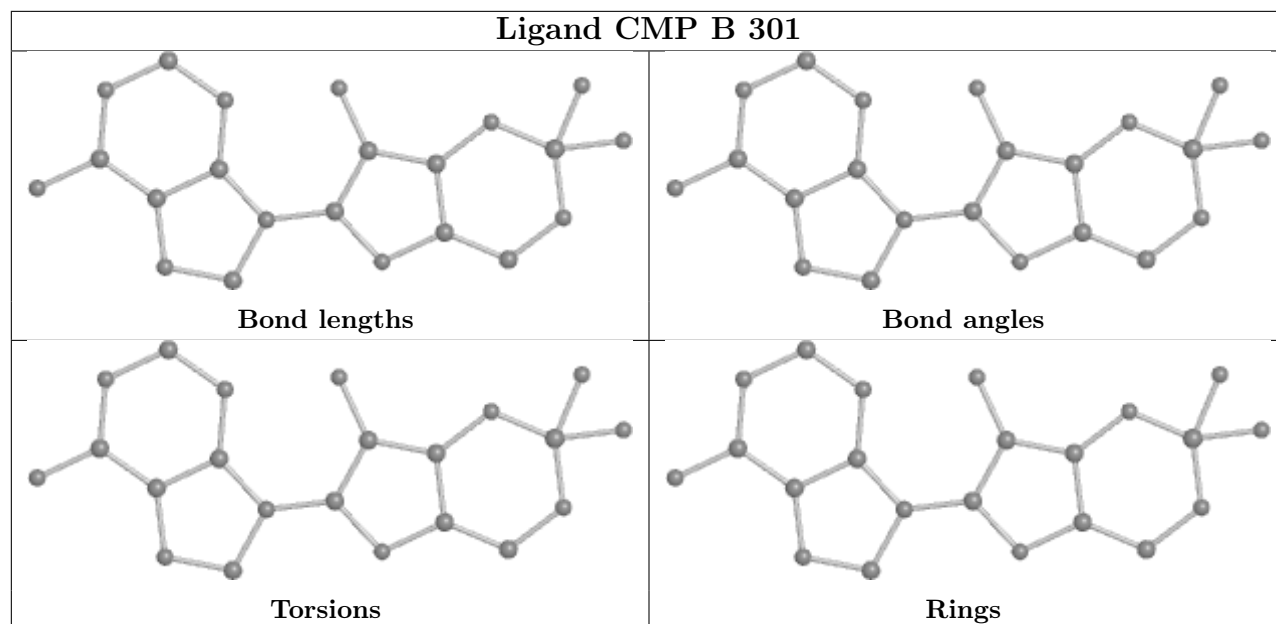
There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	206/213 (96%)	-0.04	4 (1%) 66 62	17, 34, 63, 85	0
1	B	200/213 (93%)	-0.10	3 (1%) 73 70	16, 33, 72, 88	0
All	All	406/426 (95%)	-0.07	7 (1%) 70 66	16, 34, 66, 88	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	205	GLY	6.7
1	A	162	MET	5.0
1	A	19	HIS	3.4
1	B	84	GLU	3.3
1	B	112	GLN	2.9
1	A	205	GLY	2.8
1	A	17	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](#)

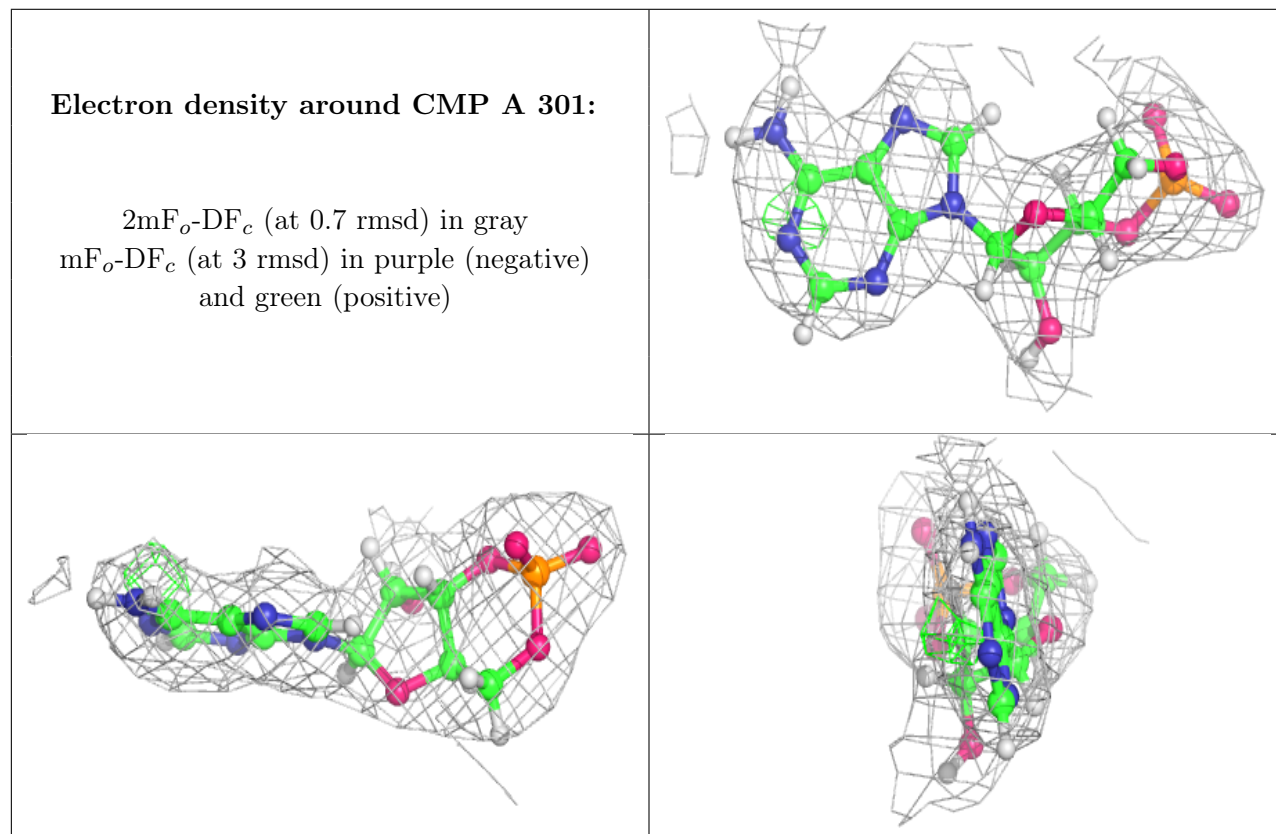
There are no monosaccharides in this entry.

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.

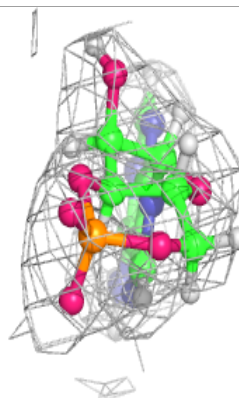
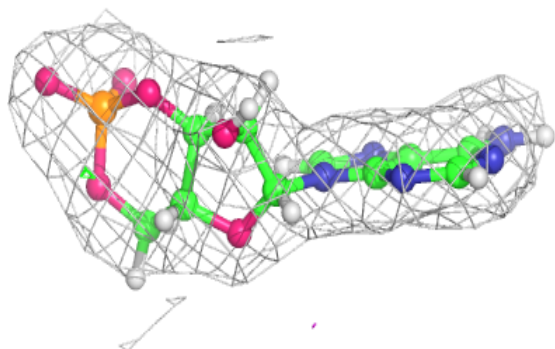
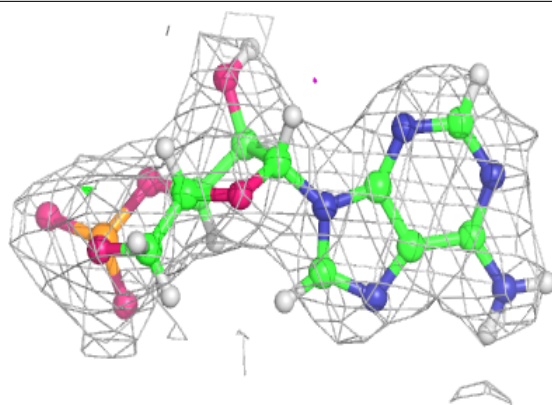
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	SO4	A	303	5/5	0.86	0.45	59,63,65,85	0
3	SO4	B	303	5/5	0.91	0.20	44,46,58,63	0
2	CMP	A	301	22/22	0.97	0.16	10,15,19,20	0
3	SO4	A	304	5/5	0.98	0.16	54,55,58,64	0
3	SO4	B	302	5/5	0.98	0.18	29,30,32,35	5
2	CMP	B	301	22/22	0.98	0.13	11,17,22,25	0
4	AU	A	305	1/1	0.98	0.07	88,88,88,88	0
3	SO4	A	302	5/5	0.99	0.14	30,31,38,43	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

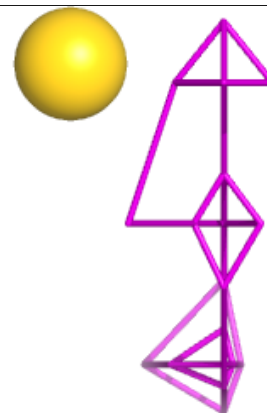
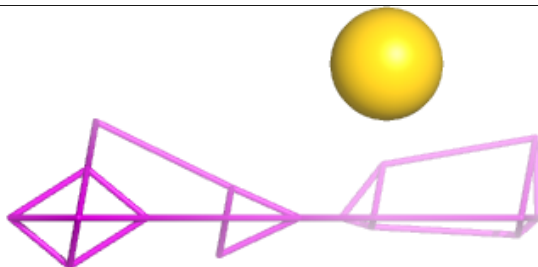
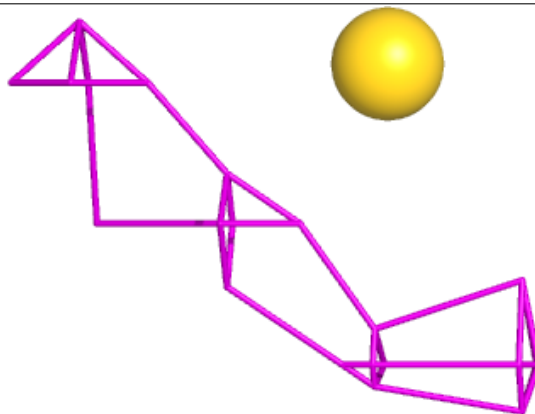


Electron density around CMP B 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around AU A 305:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.