



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 10:05 pm GMT

PDB ID : 1W4N
Title : AGAO COVALENT COMPLEX WITH TRANYLCYPROMINE
Authors : Duff, A.P.; Trambaiolo, D.M.; Langley, D.B.; Juda, G.A.; Shepard, E.M.;
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Deposited on : 2004-07-27
Resolution : 1.65 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	:	trunk28683
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

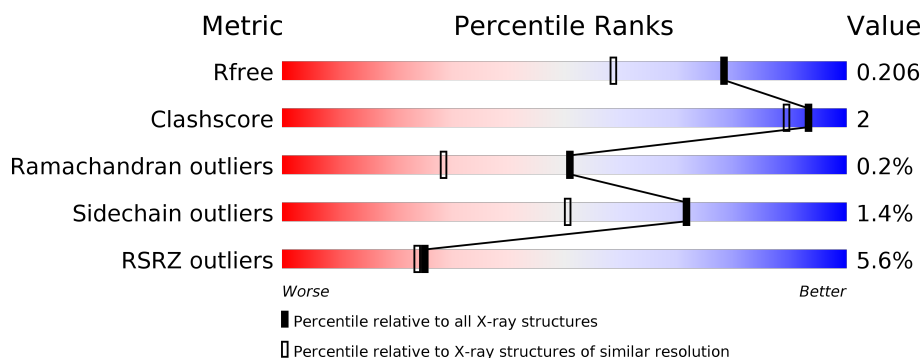
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	1368 (1.66-1.66)
Clashscore	112137	1468 (1.66-1.66)
Ramachandran outliers	110173	1438 (1.66-1.66)
Sidechain outliers	110143	1438 (1.66-1.66)
RSRZ outliers	101464	1371 (1.66-1.66)

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	646	
1	B	646	

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	SO4	A	711	-	-	-	X
4	SO4	B	711	-	-	-	X
5	GOL	A	751	-	-	-	X
5	GOL	A	752	-	-	-	X
5	GOL	A	753	-	-	-	X
5	GOL	A	755	-	-	-	X
5	GOL	B	751	-	-	-	X
5	GOL	B	752	-	-	-	X
5	GOL	B	754	-	-	-	X
5	GOL	B	755	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 19849 atoms, of which 9001 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 PHENYLETHYLAMINE OXIDASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	619	Total	C	H	N	O	S	0	2	0
			9351	3082	4478	854	927	10			
1	B	619	Total	C	H	N	O	S	0	2	0
			9351	3082	4478	854	927	10			

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

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

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Na	0	0
			1	1		
3	A	1	Total	Na	0	0
			1	1		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	0	0
			11	3	5	3		
5	A	1	Total	C	H	O	0	0
			11	3	5	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	0	0
			11	3	5	3		
5	A	1	Total	C	H	O	0	0
			11	3	5	3		
5	B	1	Total	C	H	O	0	0
			11	3	5	3		
5	B	1	Total	C	H	O	0	0
			11	3	5	3		
5	B	1	Total	C	H	O	0	0
			11	3	5	3		
5	B	1	Total	C	H	O	0	0
			11	3	5	3		

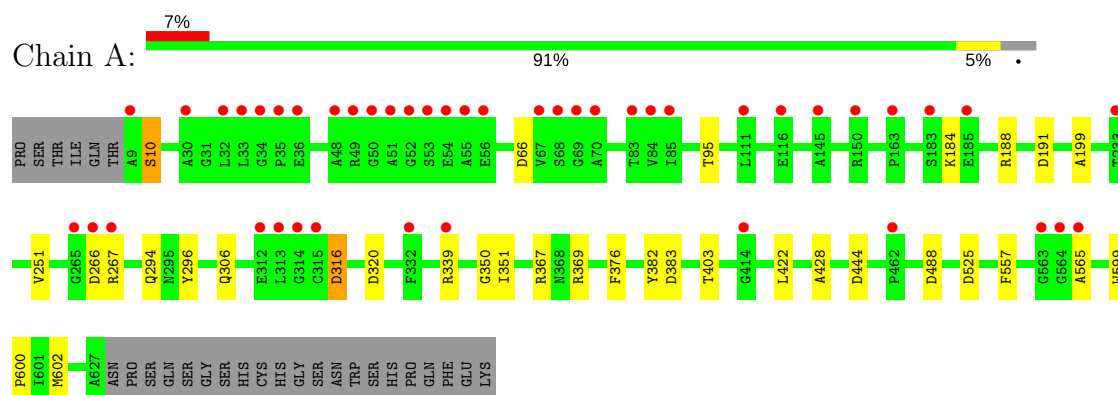
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	496	Total	O	0	0
			496	496		
6	B	538	Total	O	0	0
			538	538		

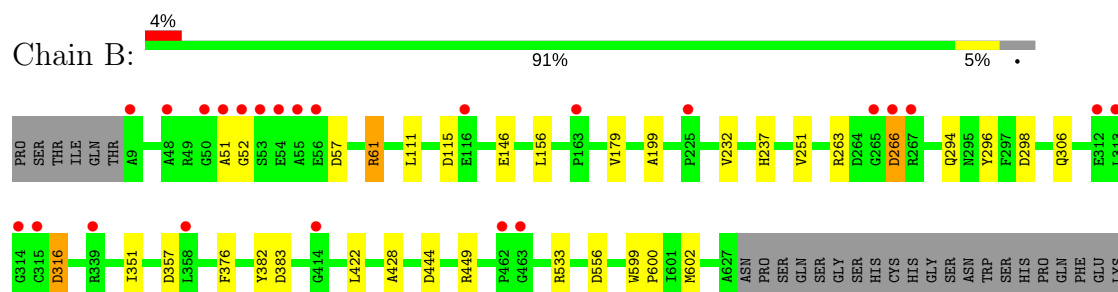
3 Residue-property plots [i](#)

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: PHENYLETHYLAMINE OXIDASE



• Molecule 1: PHENYLETHYLAMINE OXIDASE



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	158.11Å 63.00Å 184.49Å 90.00° 112.04° 90.00°	Depositor
Resolution (Å)	28.28 – 1.65 27.24 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.2 (28.28-1.65) 99.2 (27.24-1.65)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 1.65Å)	Xtrriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.171 , 0.199 0.184 , 0.206	Depositor DCC
R_{free} test set	10049 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	22.7	Xtrriage
Anisotropy	0.164	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	19849	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 76.11 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.1100e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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, TCQ, SO4, CU, NA

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	0.52	0/4983	0.80	8/6785 (0.1%)
1	B	0.60	0/4983	0.84	10/6785 (0.1%)
All	All	0.56	0/9966	0.82	18/13570 (0.1%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	61	ARG	NE-CZ-NH2	-8.83	115.89	120.30
1	B	316	ASP	CB-CG-OD2	6.58	124.22	118.30
1	B	357	ASP	CB-CG-OD2	6.46	124.11	118.30
1	A	383	ASP	CB-CG-OD2	6.35	124.01	118.30
1	A	66	ASP	CB-CG-OD2	6.28	123.95	118.30
1	B	57	ASP	CB-CG-OD2	6.08	123.78	118.30
1	B	383	ASP	CB-CG-OD2	6.02	123.72	118.30
1	A	316	ASP	CB-CG-OD2	5.91	123.62	118.30
1	B	298	ASP	CB-CG-OD2	5.68	123.41	118.30
1	B	533	ARG	NE-CZ-NH2	-5.56	117.52	120.30
1	A	191	ASP	CB-CG-OD2	5.50	123.25	118.30
1	A	320	ASP	CB-CG-OD2	5.44	123.19	118.30
1	B	449	ARG	NE-CZ-NH2	-5.43	117.58	120.30
1	A	488	ASP	CB-CG-OD2	5.43	123.19	118.30
1	A	525	ASP	CB-CG-OD2	5.18	122.97	118.30
1	B	115	ASP	CB-CG-OD2	5.12	122.91	118.30
1	B	266	ASP	CB-CG-OD2	5.04	122.83	118.30
1	A	266	ASP	CB-CG-OD2	5.01	122.81	118.30

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	4873	4478	4690	15	0
1	B	4873	4478	4689	18	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	5	0	0	1	0
4	B	5	0	0	1	0
5	A	24	20	32	1	0
5	B	30	25	40	1	0
6	A	496	0	0	2	0
6	B	538	0	0	1	0
All	All	10848	9001	9451	33	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 2.

All (33) 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:294:GLN:HG3	1:B:296:TYR:OH	1.91	0.70
4:B:711:SO4:O2	6:B:2536:HOH:O	2.11	0.67
4:A:711:SO4:O1	6:A:2490:HOH:O	2.12	0.65
1:A:565:ALA:HB1	6:A:2444:HOH:O	2.00	0.61
1:B:599:TRP:CD2	1:B:600:PRO:HA	2.35	0.61
1:B:296:TYR:CG	1:B:382:TCQ:H92	2.35	0.61
1:A:294:GLN:HG3	1:A:296:TYR:CZ	2.37	0.58
1:A:599:TRP:CD2	1:A:600:PRO:HA	2.39	0.58
1:B:251:VAL:HG22	1:B:306:GLN:HB3	1.84	0.58
1:B:61:ARG:HH22	1:B:556:ASP:CG	2.08	0.56
1:B:111:LEU:HD12	1:B:179:VAL:HG11	1.88	0.55
1:B:294:GLN:CG	1:B:296:TYR:OH	2.56	0.54
1:B:422:LEU:HD11	1:B:428:ALA:HB2	1.90	0.52
1:A:350:GLY:HA2	1:A:367:ARG:NH2	2.25	0.52
1:B:199:ALA:HA	5:B:751:GOL:H32	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:294:GLN:HG2	1:B:296:TYR:CE1	2.45	0.51
1:A:251:VAL:HG22	1:A:306:GLN:HB3	1.92	0.49
1:A:294:GLN:HG3	1:A:296:TYR:OH	2.13	0.49
1:B:61:ARG:NH2	1:B:556:ASP:OD2	2.42	0.49
1:A:422:LEU:HD11	1:A:428:ALA:HB2	1.96	0.48
1:B:296:TYR:CD2	1:B:382:TCQ:H92	2.49	0.47
1:B:294:GLN:CG	1:B:296:TYR:CZ	2.97	0.47
1:A:316:ASP:HA	1:B:351[A]:ILE:HD11	1.97	0.46
1:B:294:GLN:HG3	1:B:296:TYR:CZ	2.51	0.46
1:A:351[A]:ILE:HD11	1:B:316:ASP:HA	1.99	0.45
1:A:95:THR:O	1:A:557:PHE:HB2	2.17	0.43
1:A:296:TYR:CG	1:A:382:TCQ:H92	2.55	0.41
1:A:296:TYR:HB3	1:A:382:TCQ:H92	2.02	0.41
1:A:199:ALA:HA	5:A:751:GOL:H32	2.02	0.41
1:A:382:TCQ:H7	1:A:382:TCQ:H6	1.80	0.40
1:A:403:THR:HA	1:A:602[A]:MET:HG3	2.03	0.40
1:B:156:LEU:CD1	1:B:294:GLN:HE22	2.34	0.40
1:B:232:VAL:HA	1:B:237:HIS:O	2.22	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	618/646 (96%)	595 (96%)	22 (4%)	1 (0%)	51	28
1	B	618/646 (96%)	594 (96%)	22 (4%)	2 (0%)	44	23
All	All	1236/1292 (96%)	1189 (96%)	44 (4%)	3 (0%)	51	28

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	51	ALA
1	B	52	GLY
1	A	10	SER

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	514/537 (96%)	506 (98%)	8 (2%)	68	45
1	B	514/537 (96%)	507 (99%)	7 (1%)	71	50
All	All	1028/1074 (96%)	1013 (98%)	15 (2%)	71	48

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

Mol	Chain	Res	Type
1	A	10	SER
1	A	184	LYS
1	A	188	ARG
1	A	267	ARG
1	A	339	ARG
1	A	369	ARG
1	A	376	PHE
1	A	444	ASP
1	B	146	GLU
1	B	263	ARG
1	B	266	ASP
1	B	376	PHE
1	B	444	ASP
1	B	602[A]	MET
1	B	602[B]	MET

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	73	GLN

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Mol	Chain	Res	Type
1	A	294	GLN
1	B	294	GLN
1	B	421	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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	TCQ	A	382	1	23,25,26	1.88	6 (26%)	25,35,37	1.24	2 (8%)
1	TCQ	B	382	1	23,25,26	1.85	4 (17%)	25,35,37	1.60	4 (16%)

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	TCQ	A	382	1	-	0/10/35/37	0/2/3/3
1	TCQ	B	382	1	-	0/10/35/37	0/2/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	382	TCQ	C10-C8	-3.14	1.47	1.51
1	A	382	TCQ	C10-C8	-2.32	1.48	1.51
1	A	382	TCQ	C6-C1	2.01	1.40	1.34
1	A	382	TCQ	C9-C7	2.50	1.52	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	382	TCQ	C9-C7	2.80	1.52	1.48
1	A	382	TCQ	CA-C	2.89	1.54	1.50
1	A	382	TCQ	C3-C4	2.97	1.40	1.35
1	B	382	TCQ	C3-C4	3.42	1.41	1.35
1	B	382	TCQ	C6-C5	5.55	1.48	1.40
1	A	382	TCQ	C6-C5	5.99	1.49	1.40

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	382	TCQ	CB-CA-C	-4.54	102.67	111.41
1	A	382	TCQ	CB-CA-C	-4.30	103.12	111.41
1	B	382	TCQ	C9-C8-C10	-2.41	117.88	122.31
1	B	382	TCQ	C11-C10-C8	-2.03	117.15	121.07
1	A	382	TCQ	C6-C1-C2	2.54	120.59	118.54
1	B	382	TCQ	C6-C1-C2	4.29	122.00	118.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	382	TCQ	3	0
1	B	382	TCQ	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 4 are monoatomic - leaving 11 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	$\# Z > 2$
4	SO4	A	711	-	4,4,4	0.19	0	6,6,6	0.56	0
5	GOL	A	751	-	5,5,5	0.33	0	5,5,5	0.43	0
5	GOL	A	752	-	5,5,5	0.30	0	5,5,5	0.27	0
5	GOL	A	753	-	5,5,5	0.16	0	5,5,5	0.27	0
5	GOL	A	755	-	5,5,5	0.15	0	5,5,5	0.26	0
4	SO4	B	711	-	4,4,4	0.16	0	6,6,6	0.43	0
5	GOL	B	751	-	5,5,5	0.47	0	5,5,5	0.75	0
5	GOL	B	752	-	5,5,5	0.20	0	5,5,5	0.34	0
5	GOL	B	753	-	5,5,5	0.23	0	5,5,5	0.44	0
5	GOL	B	754	-	5,5,5	0.19	0	5,5,5	0.39	0
5	GOL	B	755	-	5,5,5	0.12	0	5,5,5	0.27	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SO4	A	711	-	-	0/0/0/0	0/0/0/0
5	GOL	A	751	-	-	0/4/4/4	0/0/0/0
5	GOL	A	752	-	-	0/4/4/4	0/0/0/0
5	GOL	A	753	-	-	0/4/4/4	0/0/0/0
5	GOL	A	755	-	-	0/4/4/4	0/0/0/0
4	SO4	B	711	-	-	0/0/0/0	0/0/0/0
5	GOL	B	751	-	-	0/4/4/4	0/0/0/0
5	GOL	B	752	-	-	0/4/4/4	0/0/0/0
5	GOL	B	753	-	-	0/4/4/4	0/0/0/0
5	GOL	B	754	-	-	0/4/4/4	0/0/0/0
5	GOL	B	755	-	-	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 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	711	SO4	1	0
5	A	751	GOL	1	0
4	B	711	SO4	1	0
5	B	751	GOL	1	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	618/646 (95%)	0.17	45 (7%)	16 15	7, 15, 39, 69	0
1	B	618/646 (95%)	-0.12	24 (3%)	40 42	4, 11, 24, 72	0
All	All	1236/1292 (95%)	0.02	69 (5%)	25 24	4, 12, 37, 72	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	53	SER	12.1
1	A	53	SER	9.5
1	A	54	GLU	8.3
1	B	52	GLY	8.3
1	A	68	SER	7.0
1	B	54	GLU	6.5
1	A	564	GLY	6.4
1	A	51	ALA	6.2
1	B	358	LEU	5.1
1	A	313	LEU	5.0
1	A	55	ALA	5.0
1	A	266	ASP	4.9
1	B	50	GLY	4.8
1	B	55	ALA	4.5
1	B	225	PRO	4.4
1	A	67	VAL	4.2
1	B	313	LEU	4.1
1	B	314	GLY	4.1
1	A	50	GLY	4.1
1	A	9	ALA	4.1
1	A	70	ALA	4.0
1	B	266	ASP	4.0
1	A	52	GLY	4.0
1	A	32	LEU	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	267	ARG	3.6
1	A	35	PRO	3.6
1	B	462	PRO	3.6
1	A	30	ALA	3.6
1	B	51	ALA	3.5
1	B	48	ALA	3.4
1	A	462	PRO	3.4
1	A	111	LEU	3.3
1	A	84	VAL	3.1
1	A	85	ILE	3.0
1	A	163	PRO	2.9
1	A	414	GLY	2.9
1	A	150	ARG	2.9
1	B	267	ARG	2.8
1	A	48	ALA	2.8
1	B	463	GLY	2.8
1	A	315	CYS	2.8
1	A	339	ARG	2.7
1	A	312	GLU	2.6
1	A	34	GLY	2.6
1	A	265	GLY	2.5
1	A	565	ALA	2.5
1	B	414	GLY	2.5
1	A	49	ARG	2.4
1	B	9	ALA	2.4
1	A	183	SER	2.4
1	B	265	GLY	2.4
1	B	315	CYS	2.4
1	A	56	GLU	2.4
1	A	185	GLU	2.4
1	B	312	GLU	2.3
1	A	145	ALA	2.3
1	B	339	ARG	2.3
1	B	163	PRO	2.3
1	A	83	THR	2.3
1	A	563	GLY	2.3
1	A	116	GLU	2.2
1	A	69	GLY	2.2
1	A	33	LEU	2.2
1	A	314	GLY	2.1
1	A	233	THR	2.1
1	A	36	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	116	GLU	2.1
1	A	332	PHE	2.1
1	B	56	GLU	2.1

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	TCQ	B	382	23/24	0.96	0.08	-	7,18,26,26	0
1	TCQ	A	382	23/24	0.94	0.10	-	10,25,42,45	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
5	GOL	B	751	6/6	0.66	0.21	11.55	39,58,59,61	0
5	GOL	A	751	6/6	0.63	0.17	10.62	49,60,62,63	0
5	GOL	B	752	6/6	0.87	0.15	7.30	50,56,58,60	0
5	GOL	A	752	6/6	0.76	0.13	6.41	57,68,70,73	0
5	GOL	B	755	6/6	0.89	0.23	3.32	37,47,53,61	0
5	GOL	A	755	6/6	0.85	0.26	3.11	45,49,56,59	0
4	SO4	A	711	5/5	0.89	0.15	2.61	49,52,60,67	0
4	SO4	B	711	5/5	0.90	0.14	2.49	55,60,69,72	0
5	GOL	A	753	6/6	0.88	0.18	2.10	35,59,65,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	GOL	B	754	6/6	0.93	0.11	2.05	24,27,29,32	0
5	GOL	B	753	6/6	0.92	0.12	0.90	34,56,58,67	0
3	NA	B	703	1/1	0.99	0.08	-0.17	26,26,26,26	0
3	NA	A	703	1/1	0.99	0.08	-0.39	32,32,32,32	0
2	CU	A	701	1/1	1.00	0.05	-2.35	29,29,29,29	0
2	CU	B	701	1/1	0.99	0.04	-2.36	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.