Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 07:07 PM BST

PDB ID	:	6VNQ
Title	:	Crystal Structure of Danio rerio Histone Deacetylase 10 in Complex with Bishydroxamic Acid Based Inhibitor
Authors	:	Herbst-Gervasoni, C.J.; Christianson, D.W.
Deposited on	:	2020-01-29
Resolution	:	2.05 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 following versions of software and data (see references ⓘ) were used in the production of this report:

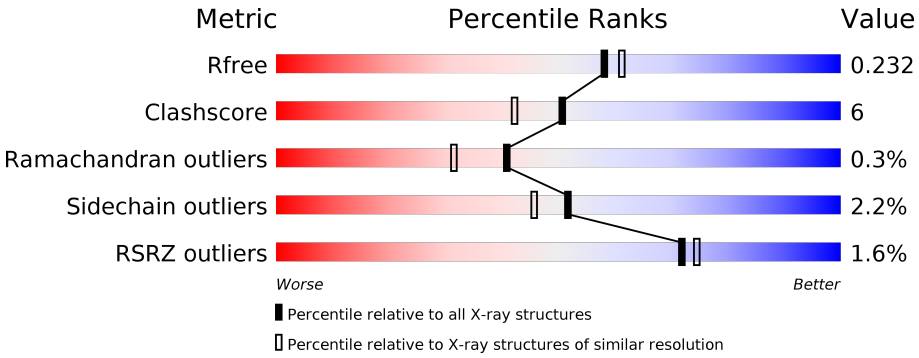
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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 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 <=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	676	% 85% 8% • 6%

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 5210 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.

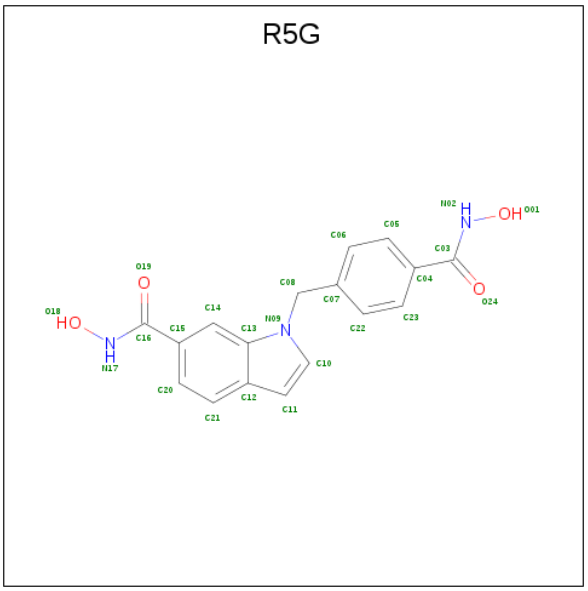
- Molecule 1 is a protein called Polyamine deacetylase HDAC10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	638	Total	C	N	O	S	0	11	0
			4880	3113	830	901	36			

There are 11 discrepancies between the modelled and reference sequences:

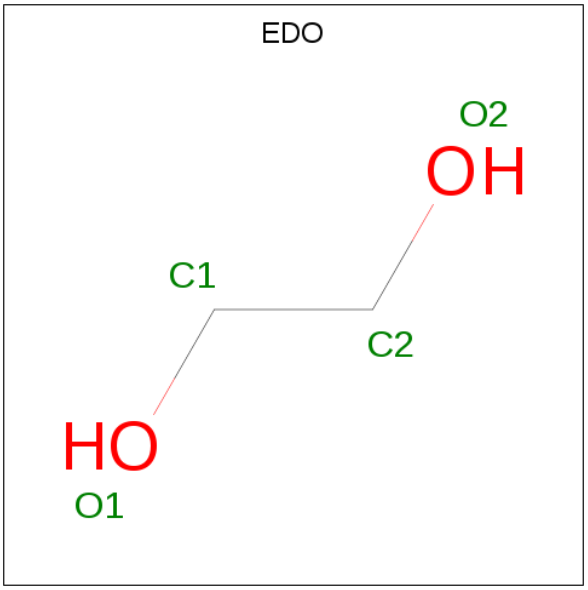
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	ALA	-	expression tag	UNP F1QCV2
A	24	GLU	ALA	engineered mutation	UNP F1QCV2
A	94	ALA	ASP	engineered mutation	UNP F1QCV2
A	154	PHE	ILE	conflict	UNP F1QCV2
A	548	THR	SER	conflict	UNP F1QCV2
A	586	GLU	GLY	conflict	UNP F1QCV2
A	593	ARG	GLY	conflict	UNP F1QCV2
A	596	ARG	THR	conflict	UNP F1QCV2
A	613	MET	THR	conflict	UNP F1QCV2
A	646	PRO	LEU	conflict	UNP F1QCV2
A	676	TRP	-	expression tag	UNP F1QCV2

- Molecule 2 is N-hydroxy-1-{[4-(hydroxycarbamoyl)phenyl]methyl}-1H-indole-6-carboxamide (three-letter code: R5G) (formula: $C_{17}H_{15}N_3O_4$) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	1
			48	34	6	8		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

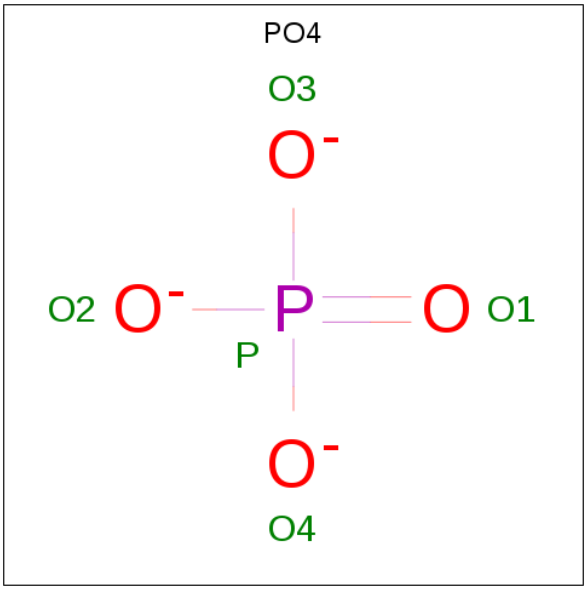
- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total K 2 2	0	0

- Molecule 6 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



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

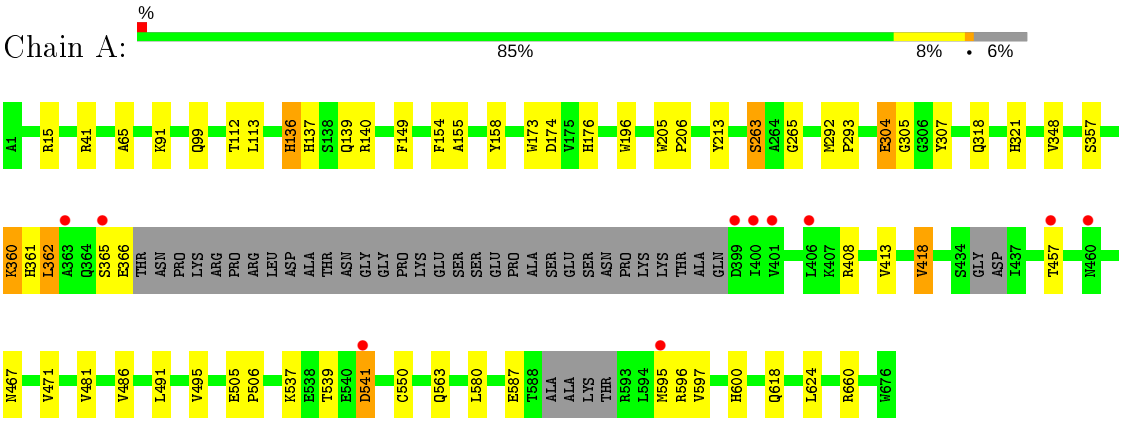
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	261	Total O 261 261	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: Polyamine deacetylase HDAC10



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	80.55Å 80.55Å 245.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	69.76 – 2.05 69.76 – 2.05	Depositor EDS
% Data completeness (in resolution range)	100.0 (69.76-2.05) 100.0 (69.76-2.05)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.71 (at 2.05Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.193 , 0.231 0.193 , 0.232	Depositor DCC
R_{free} test set	3062 reflections (5.19%)	wwPDB-VP
Wilson B-factor (Å^2)	36.1	Xtriage
Anisotropy	0.103	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 49.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.032 for -h,-k,l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	5210	wwPDB-VP
Average B, all atoms (Å^2)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.39% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: R5G, ZN, K, PO4, 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	0.46	1/5017 (0.0%)	0.57	0/6833

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	304	GLU	CD-OE2	-5.56	1.19	1.25

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	139	GLN	Mainchain

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	4880	0	4744	55	0
2	A	48	0	0	4	0
3	A	8	0	12	0	0
4	A	1	0	0	0	0
5	A	2	0	0	0	0
6	A	10	0	0	0	0
7	A	261	0	0	7	0
All	All	5210	0	4756	55	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 6.

All (55) 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:418:VAL:CG1	1:A:486:VAL:HG22	1.72	1.18
1:A:362:LEU:CG	1:A:365:SER:CB	2.22	1.16
1:A:362:LEU:HG	1:A:365:SER:CB	1.80	1.11
1:A:362:LEU:HB3	1:A:365:SER:CB	1.82	1.08
1:A:362:LEU:CB	1:A:365:SER:CB	2.32	1.08
1:A:418:VAL:HG11	1:A:486:VAL:HG22	1.44	0.98
1:A:362:LEU:HD23	1:A:365:SER:CB	2.08	0.84
1:A:362:LEU:CD2	1:A:365:SER:CB	2.56	0.84
1:A:418:VAL:HG13	1:A:486:VAL:HA	1.71	0.72
1:A:418:VAL:HG13	1:A:486:VAL:HG22	1.71	0.71
1:A:360:LYS:C	1:A:360:LYS:HD2	2.15	0.67
1:A:137:HIS:NE2	2:A:701[A]:R5G:N02	2.46	0.64
1:A:366:GLU:HG3	1:A:366:GLU:O	1.99	0.62
1:A:136:HIS:NE2	7:A:801:HOH:O	2.31	0.59
1:A:307:TYR:OH	2:A:701[A]:R5G:O24	2.20	0.58
1:A:595:MET:C	1:A:597:VAL:H	2.06	0.58
1:A:418:VAL:O	1:A:418:VAL:HG22	2.02	0.58
1:A:360:LYS:HD2	1:A:360:LYS:O	2.03	0.57
1:A:467:ASN:O	1:A:471:VAL:HG12	2.06	0.56
1:A:563:GLN:HG2	7:A:1030:HOH:O	2.06	0.55
1:A:15:ARG:NH2	7:A:812:HOH:O	2.40	0.54
1:A:587:GLU:HB2	1:A:618:GLN:HG2	1.92	0.52
1:A:505:GLU:N	1:A:506:PRO:HA	2.26	0.50
1:A:263[B]:SER:OG	1:A:304:GLU:HG3	2.10	0.50
1:A:41:ARG:HD3	7:A:971:HOH:O	2.13	0.48
1:A:65:ALA:HB3	1:A:154:PHE:HE1	1.77	0.48
1:A:418:VAL:CG1	1:A:486:VAL:CG2	2.67	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:TYR:HB2	1:A:348[B]:VAL:HG12	1.96	0.48
1:A:541:ASP:HB3	1:A:660:ARG:CZ	2.47	0.45
1:A:149:PHE:HB3	1:A:154:PHE:CZ	2.52	0.45
1:A:595:MET:HE1	1:A:624:LEU:HD21	1.99	0.45
1:A:112:THR:HG22	1:A:155:ALA:HB2	1.99	0.44
1:A:361:HIS:CD2	1:A:361:HIS:O	2.70	0.44
1:A:140[B]:ARG:NH1	7:A:807:HOH:O	2.34	0.44
1:A:176:HIS:CE1	2:A:701[A]:R5G:C23	3.01	0.44
1:A:305:GLY:HA3	2:A:701[B]:R5G:O01	2.18	0.44
1:A:292:MET:N	1:A:293:PRO:CD	2.81	0.43
1:A:99:GLN:NE2	7:A:825:HOH:O	2.52	0.43
1:A:65:ALA:HB3	1:A:154:PHE:CE1	2.53	0.43
1:A:418:VAL:HG12	1:A:486:VAL:HG22	1.87	0.43
1:A:600:HIS:HD2	7:A:970:HOH:O	2.02	0.43
1:A:360:LYS:NZ	1:A:360:LYS:O	2.35	0.42
1:A:537:LYS:O	1:A:539:THR:HG23	2.20	0.42
1:A:113:LEU:HD22	1:A:158:TYR:CG	2.55	0.42
1:A:213:TYR:CB	1:A:348[B]:VAL:HG12	2.50	0.41
1:A:205:TRP:CG	1:A:206:PRO:HA	2.55	0.41
1:A:321:HIS:HB3	1:A:408:ARG:CZ	2.50	0.41
1:A:174:ASP:OD1	1:A:265:GLY:HA3	2.21	0.41
1:A:413:VAL:O	1:A:481:VAL:HA	2.20	0.41
1:A:491:LEU:O	1:A:495:VAL:HG23	2.21	0.41
1:A:362:LEU:HB3	1:A:365:SER:CA	2.47	0.41
1:A:318:GLN:OE1	1:A:408:ARG:NH1	2.52	0.41
1:A:418:VAL:HG13	1:A:486:VAL:CG2	2.47	0.40
1:A:595:MET:HE1	1:A:624:LEU:CD2	2.51	0.40
1:A:91:LYS:HB2	1:A:91:LYS:HE3	1.90	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	641/676 (95%)	612 (96%)	27 (4%)	2 (0%)	41 31

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	357	SER
1	A	596	ARG

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	516/580 (89%)	504 (98%)	12 (2%)	50 44

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	136	HIS
1	A	173	TRP
1	A	196	TRP
1	A	263[A]	SER
1	A	263[B]	SER
1	A	360	LYS
1	A	362	LEU
1	A	418	VAL
1	A	457	THR
1	A	541	ASP
1	A	550	CYS
1	A	580	LEU

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

Mol	Chain	Res	Type
1	A	361	HIS

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

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 9 ligands modelled in this entry, 3 are monoatomic - leaving 6 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	$\#\|Z\| > 2$
6	PO4	A	708	-	4,4,4	0.95	0	6,6,6	0.60	0
3	EDO	A	702	-	3,3,3	0.52	0	2,2,2	0.15	0
2	R5G	A	701[B]	4	26,26,26	2.62	10 (38%)	33,36,36	1.62	5 (15%)
3	EDO	A	703	-	3,3,3	0.49	0	2,2,2	0.43	0
6	PO4	A	707	-	4,4,4	0.89	0	6,6,6	0.93	0
2	R5G	A	701[A]	4	26,26,26	2.27	3 (11%)	33,36,36	2.15	10 (30%)

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	702	-	-	0/1/1/1	-
2	R5G	A	701[B]	4	-	5/16/16/16	0/3/3/3
2	R5G	A	701[A]	4	-	6/16/16/16	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	703	-	-	0/1/1/1	-

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701[A]	R5G	C16-N17	7.64	1.42	1.32
2	A	701[A]	R5G	C03-N02	7.19	1.42	1.32
2	A	701[B]	R5G	C08-N09	-6.78	1.38	1.49
2	A	701[B]	R5G	C16-N17	6.14	1.40	1.32
2	A	701[B]	R5G	C04-C03	-3.94	1.42	1.50
2	A	701[B]	R5G	C13-N09	-3.70	1.34	1.39
2	A	701[B]	R5G	C03-N02	3.70	1.37	1.32
2	A	701[A]	R5G	C08-N09	-3.26	1.44	1.49
2	A	701[B]	R5G	O01-N02	-3.06	1.32	1.40
2	A	701[B]	R5G	O18-N17	-2.84	1.32	1.40
2	A	701[B]	R5G	C14-C13	-2.26	1.36	1.40
2	A	701[B]	R5G	C14-C15	-2.13	1.33	1.37
2	A	701[B]	R5G	C23-C04	-2.12	1.35	1.39

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	701[A]	R5G	C07-C08-N09	7.64	125.18	112.71
2	A	701[B]	R5G	C07-C08-N09	4.66	120.32	112.71
2	A	701[B]	R5G	C08-N09-C10	-4.03	118.80	124.73
2	A	701[A]	R5G	C08-N09-C10	-3.93	118.93	124.73
2	A	701[A]	R5G	C04-C03-N02	3.21	121.30	116.16
2	A	701[A]	R5G	C23-C04-C05	-3.10	114.17	118.59
2	A	701[A]	R5G	C06-C07-C22	-2.78	113.79	118.17
2	A	701[A]	R5G	O24-C03-N02	-2.59	118.11	122.94
2	A	701[B]	R5G	C04-C03-N02	2.56	120.26	116.16
2	A	701[A]	R5G	C11-C12-C13	2.49	108.33	106.20
2	A	701[A]	R5G	C15-C16-N17	2.33	119.90	116.16
2	A	701[B]	R5G	C06-C07-C22	-2.24	114.64	118.17
2	A	701[A]	R5G	C22-C23-C04	2.23	123.38	120.78
2	A	701[A]	R5G	C08-C07-C06	2.19	124.56	120.40
2	A	701[B]	R5G	C23-C04-C05	-2.16	115.51	118.59

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701[B]	R5G	C20-C15-C16-N17
2	A	701[A]	R5G	C07-C08-N09-C13
2	A	701[B]	R5G	C07-C08-N09-C13
2	A	701[A]	R5G	C20-C15-C16-N17
2	A	701[A]	R5G	C07-C08-N09-C10
2	A	701[B]	R5G	C20-C15-C16-O19
2	A	701[B]	R5G	C14-C15-C16-N17
2	A	701[A]	R5G	C20-C15-C16-O19
2	A	701[B]	R5G	C14-C15-C16-O19
2	A	701[A]	R5G	C14-C15-C16-N17
2	A	701[A]	R5G	C14-C15-C16-O19

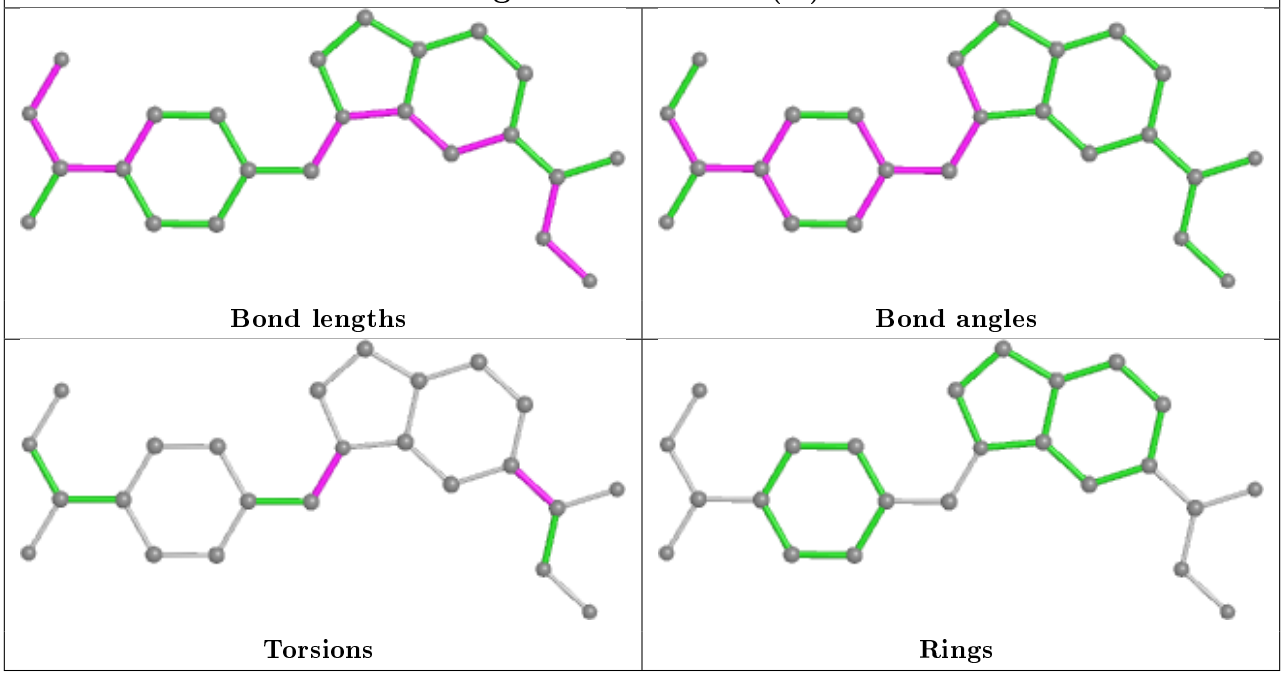
There are no ring outliers.

2 monomers are involved in 4 short contacts:

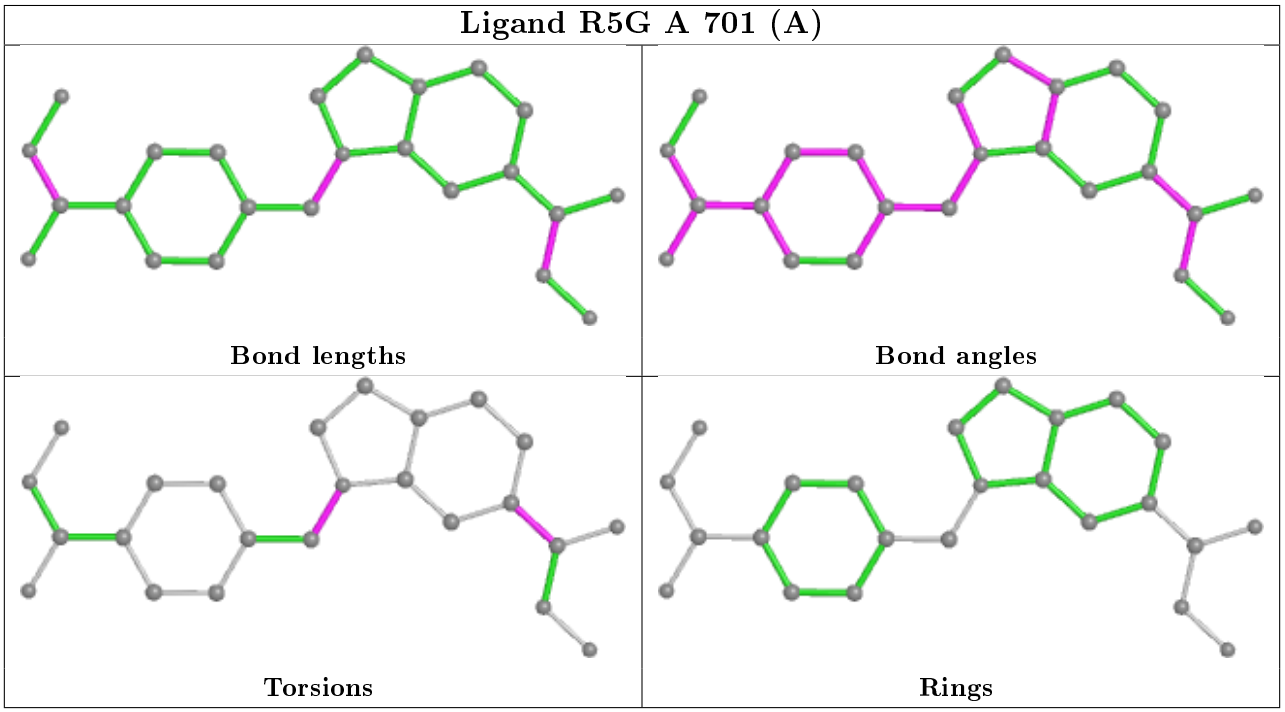
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	701[B]	R5G	1	0
2	A	701[A]	R5G	3	0

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 then 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.

Ligand R5G A 701 (B)



Ligand R5G A 701 (A)



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, 95^{th} 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	638/676 (94%)	-0.09	10 (1%) 72 74	24, 38, 75, 101	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	365	SER	4.6
1	A	363	ALA	4.1
1	A	399	ASP	3.3
1	A	457	THR	3.2
1	A	406	LEU	2.6
1	A	401	VAL	2.6
1	A	541	ASP	2.4
1	A	460	ASN	2.3
1	A	400	ILE	2.1
1	A	595	MET	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

There are no monosaccharides in this entry.

6.4 Ligands

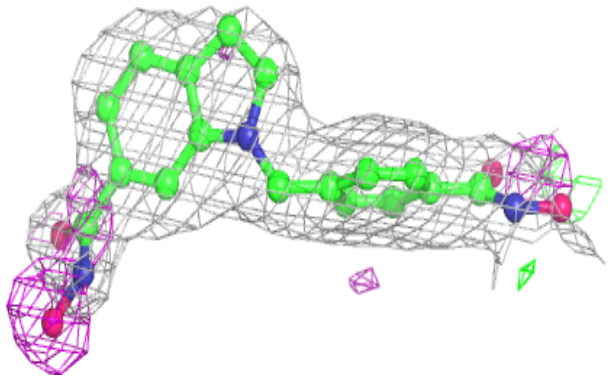
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

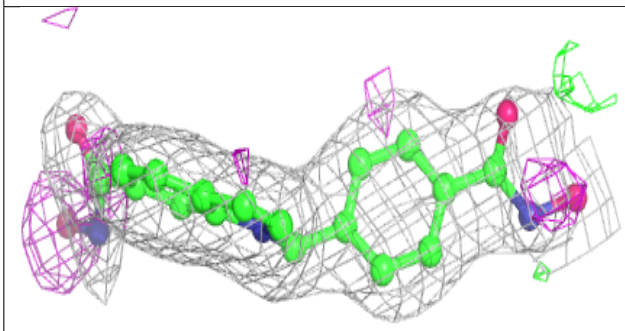
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	EDO	A	702	4/4	0.87	0.26	43,47,47,52	0
2	R5G	A	701[B]	24/24	0.88	0.17	35,42,52,64	24
2	R5G	A	701[A]	24/24	0.88	0.17	32,42,52,64	24
3	EDO	A	703	4/4	0.93	0.32	46,50,53,55	0
5	K	A	706	1/1	0.99	0.12	31,31,31,31	0
6	PO4	A	708	5/5	0.99	0.10	37,38,42,42	0
6	PO4	A	707	5/5	0.99	0.14	25,28,30,38	0
5	K	A	705	1/1	0.99	0.13	24,24,24,24	0
4	ZN	A	704	1/1	1.00	0.12	32,32,32,32	0

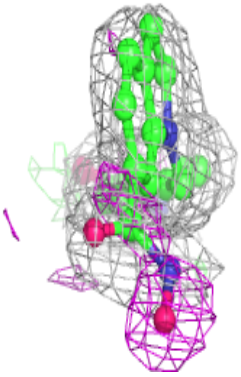
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around R5G A 701 (B):

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

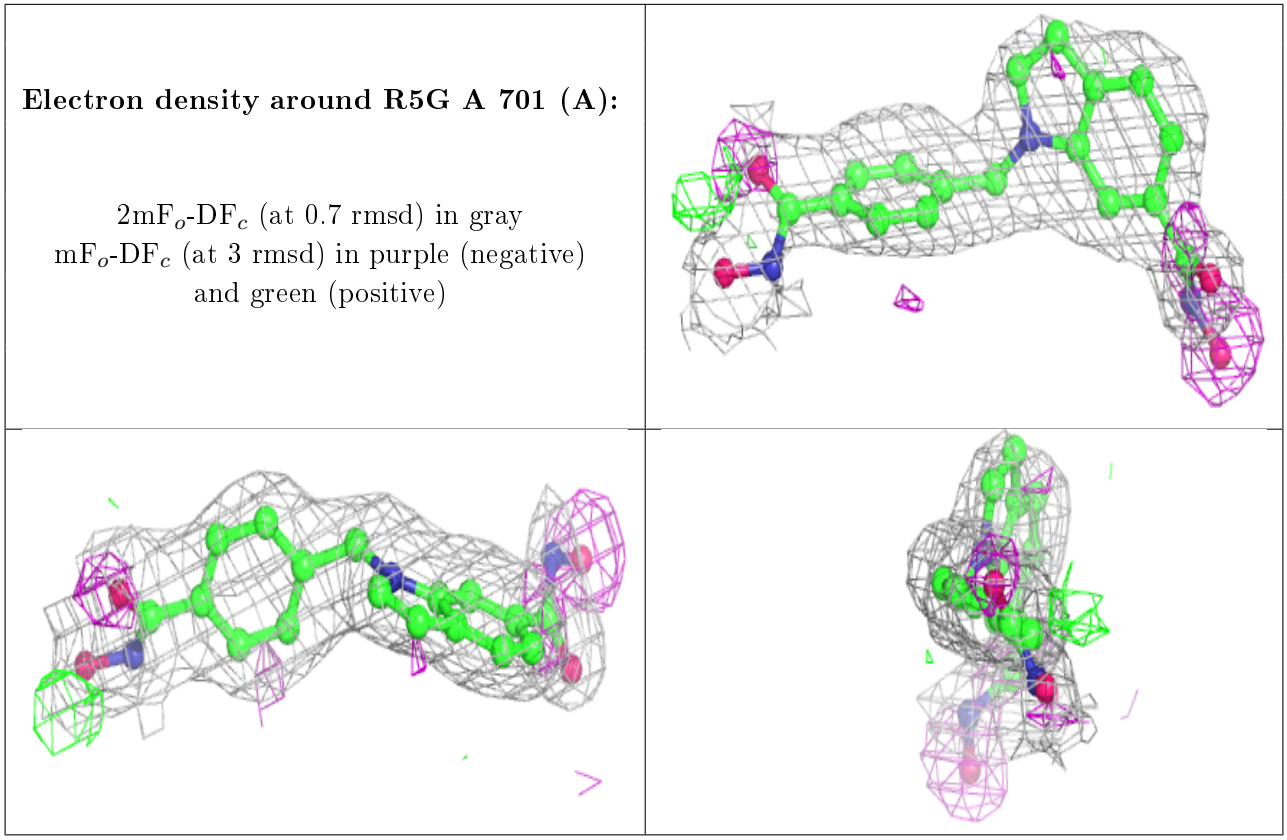






Electron density around R5G A 701 (A):

$2mF_o$-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.