



# Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2022 – 01:32 pm BST

PDB ID : 7OKZ  
Title : CRYSTAL STRUCTURE OF THE COFACTOR-DEVOID 1-H-3-HYDROXY-4- OXOQUINALDINE 2,4-DIOXYGENASE (HOD) CATALYTICALLY INACTIVE H251A VARIANT COMPLEXED WITH 2-METHYL- QUINOLIN-4(1H)-ONE UNDER HYPEROXIC CONDITIONS  
Authors : Bui, S.; Steiner, R.A.  
Deposited on : 2021-05-18  
Resolution : 2.10 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.28.1  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0267  
CCP4 : 7.1.010 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.28.1

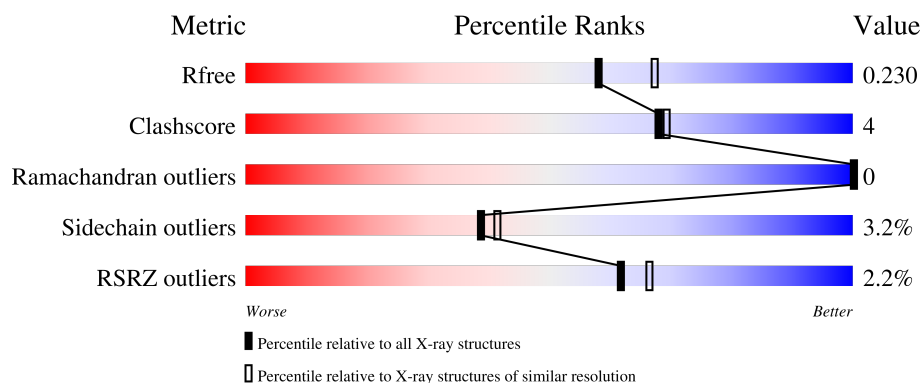
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

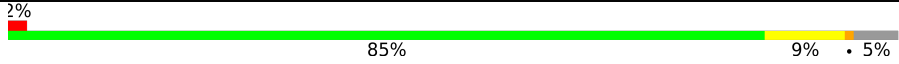

The reported resolution of this entry is 2.10 Å.

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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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  $\geq 3$ , 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	AAA	288	
1	BBB	288	

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
5	GOL	BBB	301	-	-	X	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4867 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 1H-3-hydroxy-4-oxoquinoline 2,4-dioxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	274	Total	C	N	O	S	0	6	0
			2282	1453	401	420	8			
1	BBB	274	Total	C	N	O	S	0	4	0
			2267	1446	398	415	8			

There are 28 discrepancies between the modelled and reference sequences:

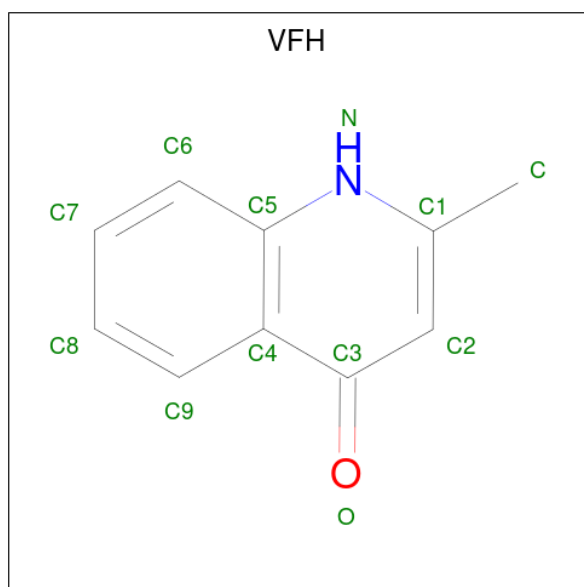
Chain	Residue	Modelled	Actual	Comment	Reference
AAA	-11	MET	-	initiating methionine	UNP O31266
AAA	-10	ARG	-	expression tag	UNP O31266
AAA	-9	GLY	-	expression tag	UNP O31266
AAA	-8	SER	-	expression tag	UNP O31266
AAA	-7	HIS	-	expression tag	UNP O31266
AAA	-6	HIS	-	expression tag	UNP O31266
AAA	-5	HIS	-	expression tag	UNP O31266
AAA	-4	HIS	-	expression tag	UNP O31266
AAA	-3	HIS	-	expression tag	UNP O31266
AAA	-2	HIS	-	expression tag	UNP O31266
AAA	-1	GLY	-	expression tag	UNP O31266
AAA	0	SER	-	expression tag	UNP O31266
AAA	69	SER	CYS	engineered mutation	UNP O31266
AAA	251	ALA	HIS	engineered mutation	UNP O31266
BBB	-11	MET	-	initiating methionine	UNP O31266
BBB	-10	ARG	-	expression tag	UNP O31266
BBB	-9	GLY	-	expression tag	UNP O31266
BBB	-8	SER	-	expression tag	UNP O31266
BBB	-7	HIS	-	expression tag	UNP O31266
BBB	-6	HIS	-	expression tag	UNP O31266
BBB	-5	HIS	-	expression tag	UNP O31266
BBB	-4	HIS	-	expression tag	UNP O31266
BBB	-3	HIS	-	expression tag	UNP O31266
BBB	-2	HIS	-	expression tag	UNP O31266
BBB	-1	GLY	-	expression tag	UNP O31266

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Chain	Residue	Modelled	Actual	Comment	Reference
BBB	0	SER	-	expression tag	UNP O31266
BBB	69	SER	CYS	engineered mutation	UNP O31266
BBB	251	ALA	HIS	engineered mutation	UNP O31266

- Molecule 2 is 2-methyl-quinolin-4(1H)-one (three-letter code: VFH) (formula:  $C_{10}H_9NO$ ) (labeled as "Ligand of Interest" by depositor).

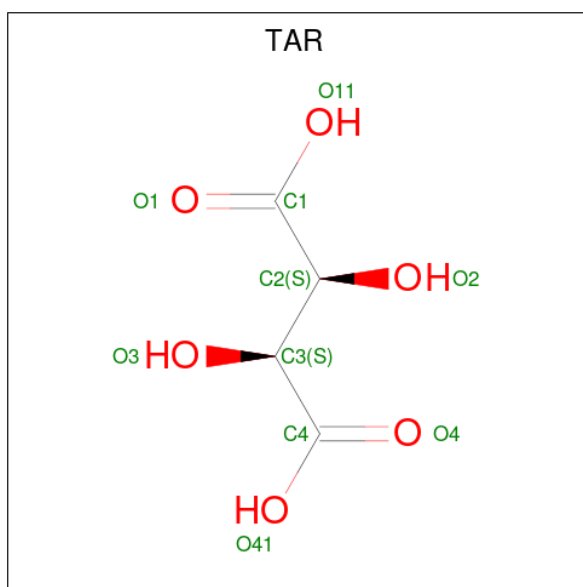


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	AAA	1	Total	C	N	O	0	0
			12	10	1	1		
2	BBB	1	Total	C	N	O	0	0
			12	10	1	1		

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	AAA	1	Total	K	0	0
			1	1		
3	BBB	1	Total	K	0	0
			1	1		

- Molecule 4 is D(-)-TARTARIC ACID (three-letter code: TAR) (formula:  $C_4H_6O_6$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	AAA	1	Total	C	O	0	0
			10	4	6		

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	AAA	1	Total	C	O	0	0
			6	3	3		
5	AAA	1	Total	C	O	0	1
			12	6	6		
5	BBB	1	Total	C	O	0	0
			6	3	3		

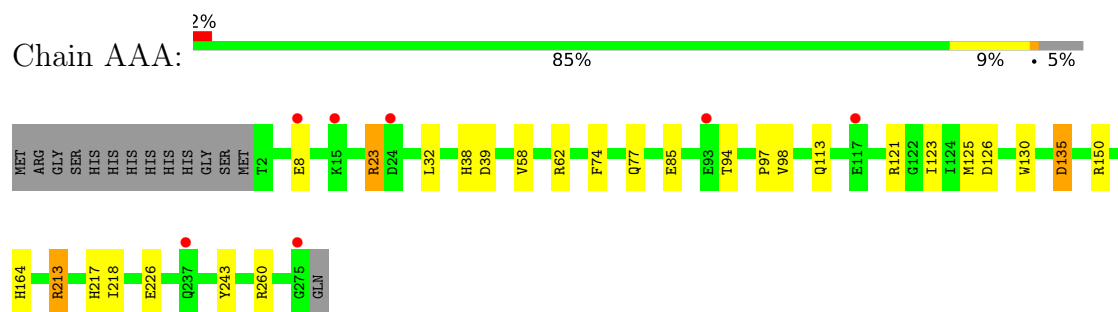
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	AAA	132	Total 135	O 135	0	3
6	BBB	123	Total 123	O 123	0	0

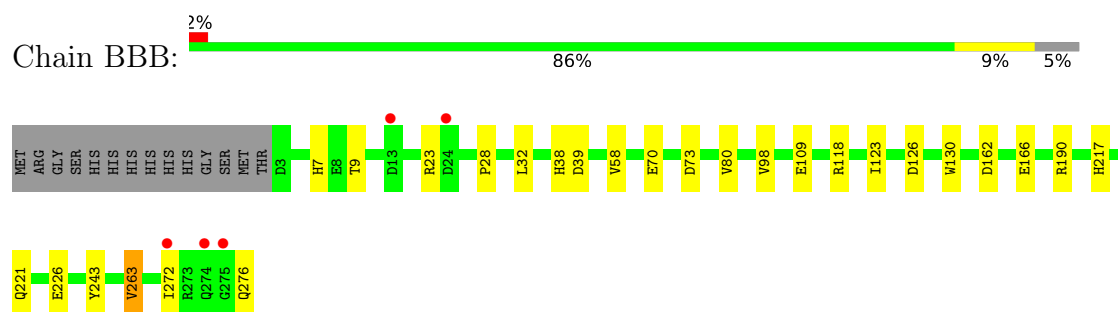
### 3 Residue-property plots [i](#)

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: 1H-3-hydroxy-4-oxoquinaldine 2,4-dioxygenase



- Molecule 1: 1H-3-hydroxy-4-oxoquinaldine 2,4-dioxygenase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	120.52Å 120.52Å 44.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.26 – 2.10 60.26 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.4 (60.26-2.10) 99.4 (60.26-2.10)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.47 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, $R_{free}$	0.190 , 0.228 0.194 , 0.230	Depositor DCC
$R_{free}$ test set	1793 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	34.8	Xtriage
Anisotropy	0.024	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	(Not available) , (Not available)	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.39$ , $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.083 for h,-k,-l	Xtriage
Reported twinning fraction	0.881 for H, K, L 0.119 for -K, -H, -L	Depositor
Outliers	0 of 37750 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4867	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.12% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: K, TAR, VFH, GOL

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	AAA	0.65	0/2352	0.77	3/3200 (0.1%)
1	BBB	0.65	0/2337	0.74	0/3179
All	All	0.65	0/4689	0.75	3/6379 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AAA	260	ARG	NE-CZ-NH2	-6.73	116.93	120.30
1	AAA	260	ARG	NE-CZ-NH1	5.36	122.98	120.30
1	AAA	213	ARG	CB-CA-C	-5.11	100.19	110.40

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	AAA	2282	0	2170	16	0
1	BBB	2267	0	2162	20	0
2	AAA	12	0	0	0	0
2	BBB	12	0	0	0	0
3	AAA	1	0	0	0	0
3	BBB	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	AAA	10	0	4	1	0
5	AAA	18	0	24	0	0
5	BBB	6	0	8	7	0
6	AAA	135	0	0	6	0
6	BBB	123	0	0	4	1
All	All	4867	0	4368	34	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 4.

All (34) 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:BBB:263[B]:VAL:CG1	5:BBB:301:GOL:H12	2.07	0.85
1:BBB:263[B]:VAL:HG11	5:BBB:301:GOL:H12	1.68	0.74
1:BBB:263[B]:VAL:HG11	5:BBB:301:GOL:C1	2.27	0.65
1:AAA:135[A]:ASP:CG	6:AAA:407:HOH:O	2.35	0.64
1:BBB:263[B]:VAL:CG1	5:BBB:301:GOL:C1	2.75	0.62
1:BBB:263[A]:VAL:HG21	5:BBB:301:GOL:H2	1.86	0.58
1:AAA:32:LEU:HB2	1:AAA:58:VAL:HG22	1.86	0.58
1:AAA:150:ARG:NH1	6:AAA:410:HOH:O	2.40	0.55
1:AAA:217:HIS:O	1:AAA:243:TYR:HA	2.07	0.54
1:BBB:32:LEU:HB2	1:BBB:58:VAL:HG22	1.90	0.53
1:BBB:162:ASP:HB2	1:BBB:221[B]:GLN:NE2	2.24	0.53
1:BBB:217:HIS:O	1:BBB:243:TYR:HA	2.09	0.52
1:AAA:85:GLU:HA	6:AAA:422:HOH:O	2.11	0.49
1:BBB:263[B]:VAL:HG11	5:BBB:301:GOL:C2	2.42	0.49
1:AAA:8:GLU:O	6:AAA:401:HOH:O	2.20	0.48
1:BBB:28:PRO:HG3	1:BBB:272[B]:ILE:HG21	1.96	0.48
1:BBB:166:GLU:OE2	6:BBB:401:HOH:O	2.20	0.47
1:AAA:94:THR:HG22	6:AAA:498:HOH:O	2.15	0.47
1:BBB:9:THR:HB	6:BBB:416:HOH:O	2.16	0.46
1:AAA:77:GLN:HB2	1:BBB:190[B]:ARG:HD2	1.99	0.45
1:BBB:118:ARG:NH2	6:BBB:408:HOH:O	2.37	0.45
1:BBB:263[B]:VAL:HG12	5:BBB:301:GOL:H12	1.97	0.43
1:BBB:7:HIS:CD2	6:BBB:476:HOH:O	2.72	0.42
1:AAA:98:VAL:HA	1:AAA:123:ILE:O	2.20	0.42
1:BBB:226:GLU:H	1:BBB:226:GLU:CD	2.24	0.42
1:AAA:226:GLU:H	1:AAA:226:GLU:CD	2.23	0.41
1:AAA:62:ARG:HB2	1:AAA:74:PHE:CE1	2.55	0.41
1:AAA:164:HIS:HD2	4:AAA:303:TAR:O11	2.04	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BBB:80:VAL:HG21	1:BBB:109:GLU:HB3	2.03	0.41
1:BBB:98:VAL:HA	1:BBB:123:ILE:O	2.19	0.41
1:AAA:113:GLN:NE2	1:BBB:73:ASP:HB2	2.36	0.41
1:AAA:97:PRO:HD2	1:AAA:121:ARG:O	2.21	0.41
1:AAA:125:MET:HA	1:AAA:218:ILE:O	2.20	0.41
1:AAA:23:ARG:NH1	6:AAA:424:HOH:O	2.53	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)
6:BBB:419:HOH:O	6:BBB:496:HOH:O[4_554]	2.16	0.04

## 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	AAA	278/288 (96%)	271 (98%)	7 (2%)	0	100	100
1	BBB	276/288 (96%)	268 (97%)	8 (3%)	0	100	100
All	All	554/576 (96%)	539 (97%)	15 (3%)	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	AAA	241/247 (98%)	233 (97%)	8 (3%)	38	40
1	BBB	239/247 (97%)	230 (96%)	9 (4%)	33	34
All	All	480/494 (97%)	463 (96%)	17 (4%)	39	38

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	23	ARG
1	AAA	38	HIS
1	AAA	39	ASP
1	AAA	126	ASP
1	AAA	130	TRP
1	AAA	135[A]	ASP
1	AAA	135[B]	ASP
1	AAA	213	ARG
1	BBB	23	ARG
1	BBB	38	HIS
1	BBB	39	ASP
1	BBB	70	GLU
1	BBB	126	ASP
1	BBB	130	TRP
1	BBB	263[A]	VAL
1	BBB	263[B]	VAL
1	BBB	276	GLN

Sometimes 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 ⓘ

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, 2 are monoatomic - leaving 7 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	$\# Z  > 2$
2	VFH	BBB	302	-	12,13,13	1.56	2 (16%)	15,18,18	2.72	3 (20%)
5	GOL	AAA	305[B]	-	5,5,5	0.11	0	5,5,5	0.23	0
5	GOL	AAA	305[A]	-	5,5,5	0.09	0	5,5,5	0.24	0
5	GOL	BBB	301	-	5,5,5	0.12	0	5,5,5	0.40	0
4	TAR	AAA	303	-	3,9,9	0.12	0	6,12,12	1.58	1 (16%)
2	VFH	AAA	301	-	12,13,13	1.54	2 (16%)	15,18,18	2.80	3 (20%)
5	GOL	AAA	304	3	5,5,5	0.13	0	5,5,5	0.43	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 Torsions and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	VFH	BBB	302	-	-	-	0/2/2/2
5	GOL	AAA	305[B]	-	-	1/4/4/4	-
5	GOL	AAA	305[A]	-	-	2/4/4/4	-
5	GOL	BBB	301	-	-	1/4/4/4	-
4	TAR	AAA	303	-	-	1/4/12/12	-
2	VFH	AAA	301	-	-	-	0/2/2/2
5	GOL	AAA	304	3	-	2/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	AAA	301	VFH	C2-C3	3.95	1.45	1.37
2	BBB	302	VFH	C2-C3	3.61	1.44	1.37
2	BBB	302	VFH	C1-N	2.54	1.38	1.33
2	AAA	301	VFH	C1-N	2.44	1.38	1.33

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	AAA	301	VFH	C4-C5-N	-8.15	119.18	123.60
2	BBB	302	VFH	C4-C5-N	-7.79	119.37	123.60
2	BBB	302	VFH	C4-C3-C2	-6.22	116.16	123.05
2	AAA	301	VFH	C4-C3-C2	-6.17	116.22	123.05
4	AAA	303	TAR	C1-C2-C3	-3.15	106.32	113.11
2	AAA	301	VFH	C1-N-C5	2.52	123.71	118.40
2	BBB	302	VFH	C1-N-C5	2.25	123.14	118.40

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	AAA	304	GOL	C1-C2-C3-O3
5	AAA	305[A]	GOL	O1-C1-C2-C3
5	AAA	304	GOL	O2-C2-C3-O3
5	AAA	305[A]	GOL	O1-C1-C2-O2
5	BBB	301	GOL	O2-C2-C3-O3
4	AAA	303	TAR	O2-C2-C3-O3
5	AAA	305[B]	GOL	C1-C2-C3-O3

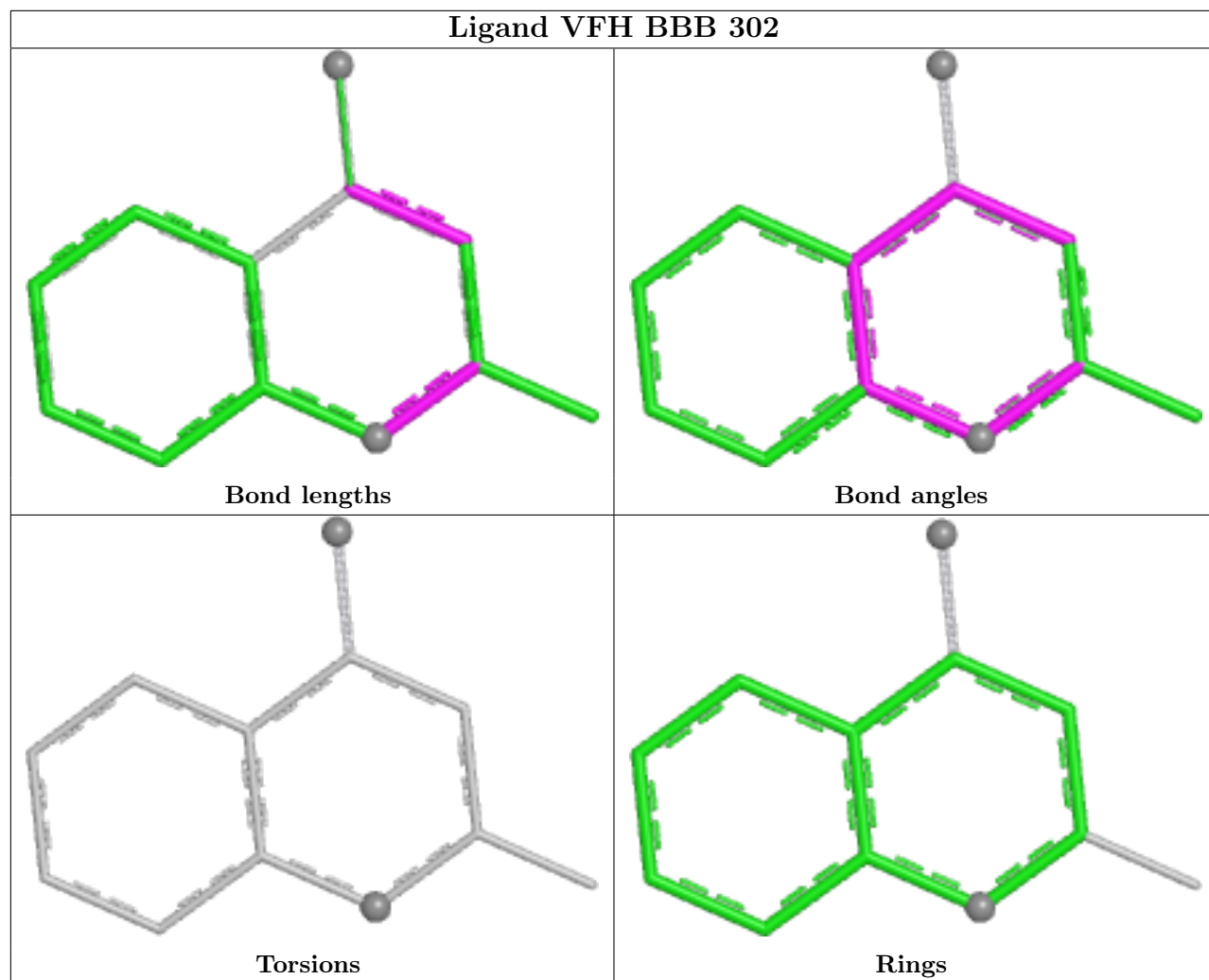
There are no ring outliers.

2 monomers are involved in 8 short contacts:

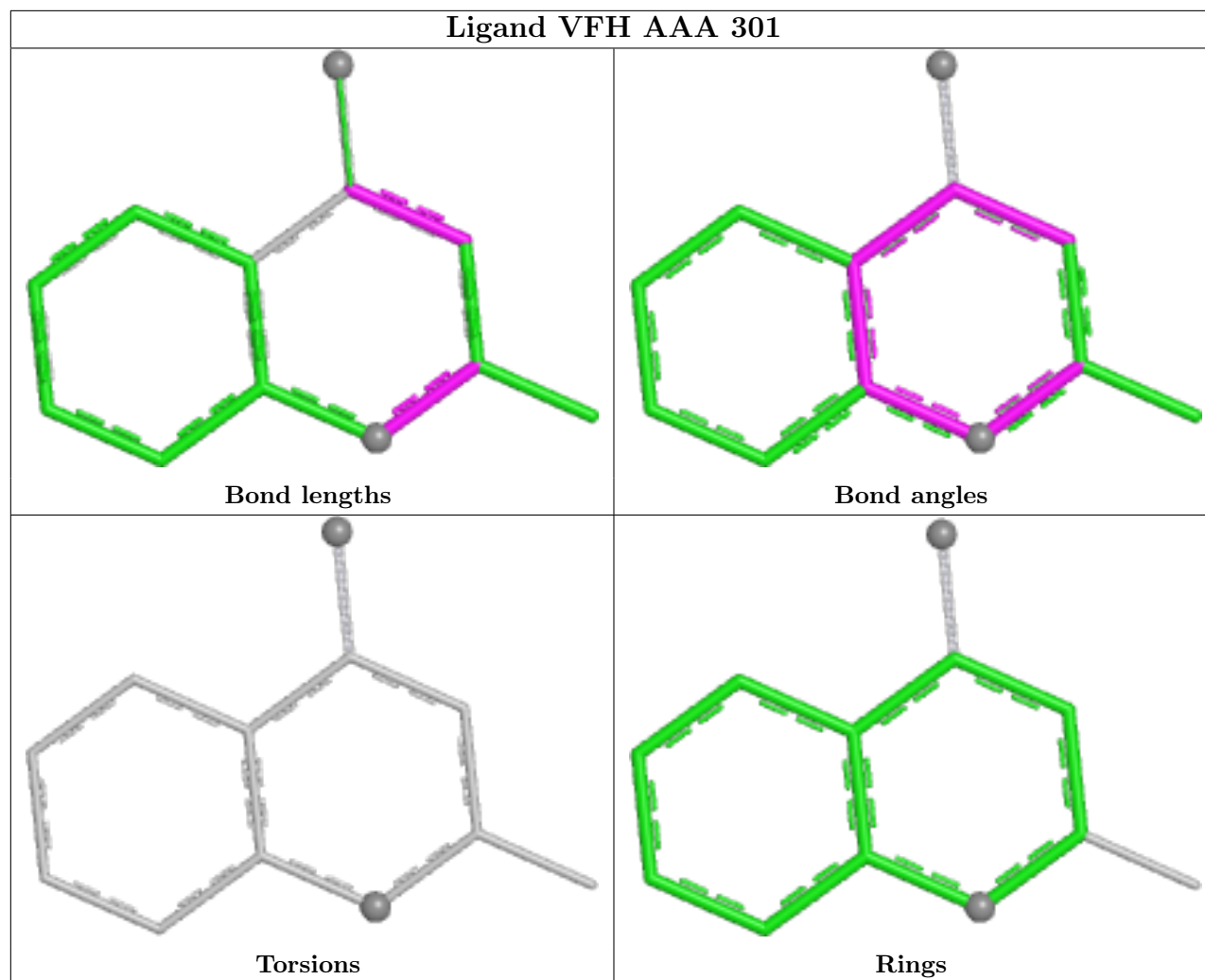
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	BBB	301	GOL	7	0
4	AAA	303	TAR	1	0

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 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.

## Ligand VFH BBB 302







## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	AAA	274/288 (95%)	0.20	7 (2%) 56 61	18, 35, 54, 73	0
1	BBB	274/288 (95%)	0.22	5 (1%) 68 72	25, 37, 57, 80	0
All	All	548/576 (95%)	0.21	12 (2%) 62 66	18, 36, 56, 80	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	24	ASP	4.8
1	AAA	93	GLU	3.4
1	BBB	13	ASP	3.3
1	AAA	275	GLY	3.3
1	BBB	275	GLY	2.9
1	BBB	24	ASP	2.7
1	AAA	237	GLN	2.7
1	BBB	274	GLN	2.5
1	AAA	15	LYS	2.3
1	AAA	117	GLU	2.2
1	BBB	272[A]	ILE	2.2
1	AAA	8	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 monosaccharides in this entry.

## 6.4 Ligands

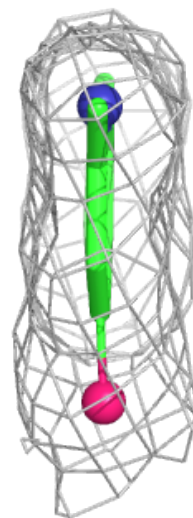
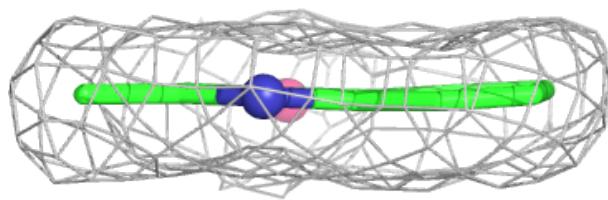
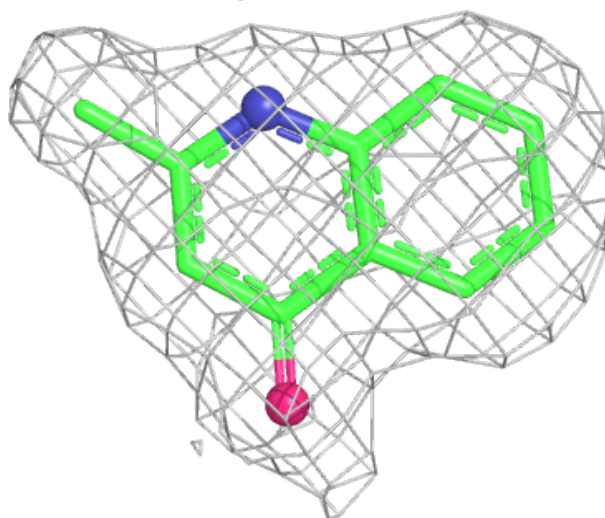
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	GOL	AAA	304	6/6	0.77	0.32	50,55,59,71	0
4	TAR	AAA	303	10/10	0.82	0.25	34,43,49,50	0
5	GOL	BBB	301	6/6	0.86	0.23	33,38,40,40	0
5	GOL	AAA	305[B]	6/6	0.88	0.20	29,32,36,39	6
5	GOL	AAA	305[A]	6/6	0.88	0.20	38,42,45,47	6
2	VFH	AAA	301	12/12	0.92	0.10	26,26,28,29	0
2	VFH	BBB	302	12/12	0.94	0.11	26,28,29,29	0
3	K	BBB	303	1/1	0.97	0.08	45,45,45,45	0
3	K	AAA	302	1/1	1.00	0.11	39,39,39,39	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.

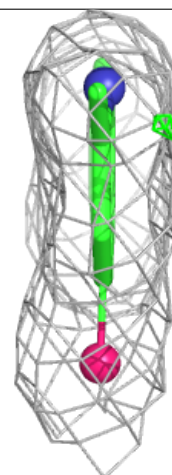
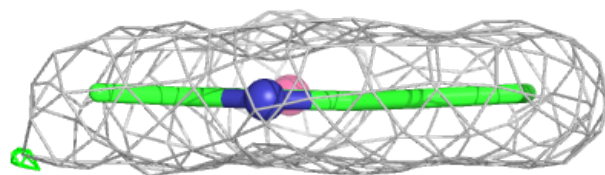
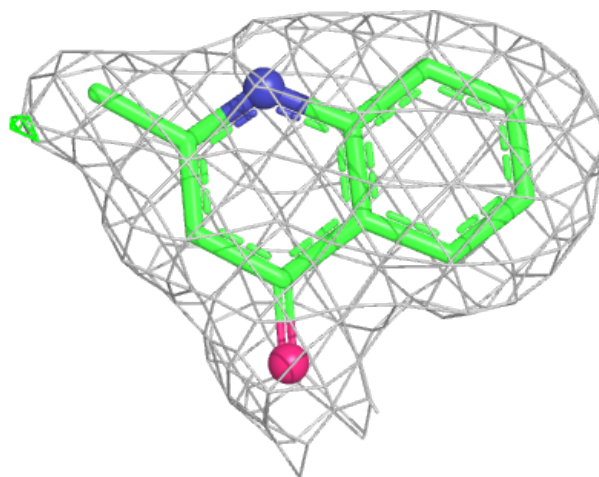
**Electron density around VFH AAA 301:**

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



**Electron density around VFH BBB 302:**

$2mF_o-DF_c$  (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.