

Full wwPDB X-ray Structure Validation Report (i)

Mar 25, 2024 – 06:07 PM EDT

PDB ID	:	8VHL
Title	:	Structure of DHODH in Complex with Ligand 17
Authors	:	Shaffer, P.L.
Deposited on		
Resolution	:	1.93 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 (i) 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 (1)) were used in the production of this report:

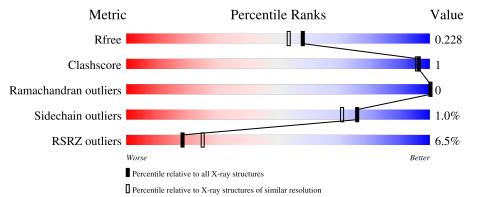
Xtriage (Phenix) EDS buster-report Percentile statistics	: : :	4.02b-467 1.8.5 (274361), CSD as541be (2020) 1.13 2.36.1 1.1.7 (2018) 20191225.v01 (using entries in the PDB archive December 25th 2019) 5.8.0158
-		
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac		
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 1.93 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	4310 (1.96-1.92)
Clashscore	141614	1023 (1.94-1.94)
Ramachandran outliers	138981	1007 (1.94-1.94)
Sidechain outliers	138945	1007 (1.94-1.94)
RSRZ outliers	127900	4250 (1.96-1.92)

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 <=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	
			6%	
1	А	369	90%	5%••



$8 \mathrm{VHL}$

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 3225 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.

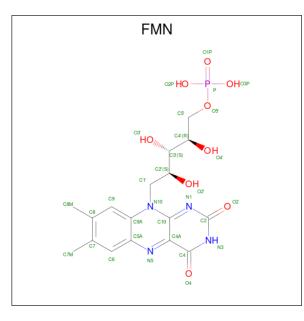
• Molecule 1 is a protein called Dihydroorotate dehydrogenase (quinone), mitochondrial.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	А	355	Total	С	Ν	0	\mathbf{S}	22	11	0
	11	000	2807	1762	514	527	4		11	U

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	28	GLY	-	expression tag	UNP Q02127
А	29	ALA	-	expression tag	UNP Q02127

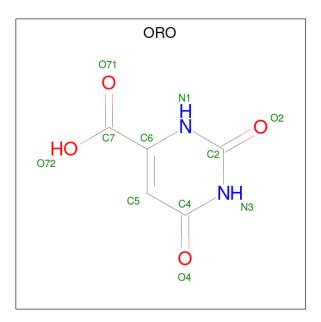
• Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: C₁₇H₂₁N₄O₉P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	А	1	Total	C	Ν	0	Р	0	0
			31	17	4	9	T		

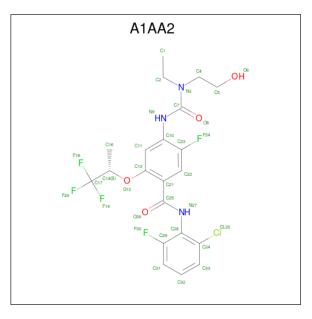
• Molecule 3 is OROTIC ACID (three-letter code: ORO) (formula: $C_5H_4N_2O_4$).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	А	1	Total 11	С 5	N 2	0 4	0	0

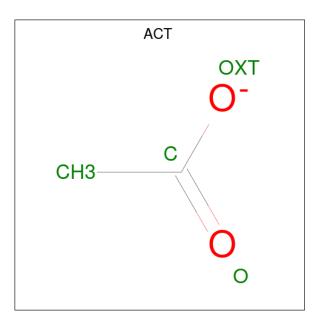
• Molecule 4 is N-(2-chloro-6-fluorophenyl)-4-[ethyl(2-hydroxyethyl)carbamamido]-5-fluoro -2-{[(2S)-1,1,1-trifluoropropan-2-yl]oxy}benzamide (three-letter code: A1AA2) (formula: $C_{21}H_{21}ClF_5N_3O_4$) (labeled as "Ligand of Interest" by depositor).



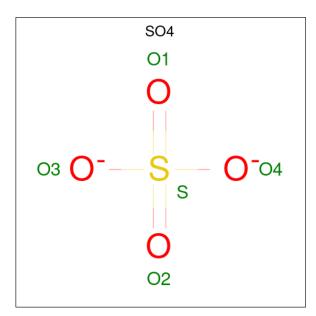
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
4	А	1	Total 34	C 21		F 5	N 3	0 4	0	0

• Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).





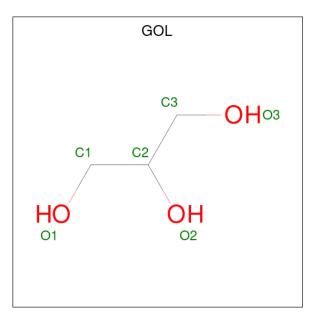
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
5	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
5	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0



Mol	Chain	Residues	Atc	\mathbf{ms}		ZeroOcc	AltConf
6	А	1	Total 5	0 4	S 1	0	0

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0

• Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0
7	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0
7	А	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 6 3 3 \end{array}$	0	0
7	А	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 6 3 3 \end{array}$	0	0

• Molecule 8 is water.

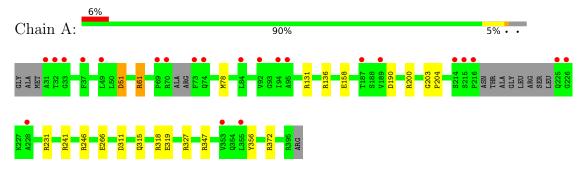
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	А	289	Total O 291 291	0	2



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: Dihydroorotate dehydrogenase (quinone), mitochondrial





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants	90.74Å 90.74 Å 122.71 Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	78.59 - 1.93	Depositor
Resolution (A)	39.29 - 1.93	EDS
% Data completeness	$95.6\ (78.59-1.93)$	Depositor
(in resolution range)	95.6 (39.29-1.93)	EDS
R _{merge}	0.07	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.54 (at 1.94 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
B B.	0.168 , 0.199	Depositor
R, R_{free}	0.198 , 0.228	DCC
R_{free} test set	2042 reflections $(4.80%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	28.3	Xtriage
Anisotropy	0.183	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.36, 53.0	EDS
L-test for twinning ²	$< L > = 0.49, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3225	wwPDB-VP
Average B, all atoms $(Å^2)$	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.37% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: ORO, A1AA2, SO4, GOL, ACT, FMN

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).

Mal	Chain	Bo	nd lengths	Bo	ond angles
Mol Chain		RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.97	3/2852~(0.1%)	1.07	18/3851~(0.5%)

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	А	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\operatorname{Observed}(\operatorname{\AA})$	$\mathrm{Ideal}(\mathrm{\AA})$
1	А	61[A]	ARG	C-O	6.44	1.35	1.23
1	А	61[B]	ARG	C-O	6.44	1.35	1.23
1	А	158	GLU	CD-OE2	-5.10	1.20	1.25

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	51	ASP	CB-CG-OD1	-7.47	111.57	118.30
1	А	347	ARG	NE-CZ-NH1	7.40	124.00	120.30
1	А	136	ARG	NE-CZ-NH1	7.34	123.97	120.30
1	А	231	ARG	NE-CZ-NH1	7.16	123.88	120.30
1	А	347	ARG	NE-CZ-NH2	-7.16	116.72	120.30
1	А	327	ARG	NE-CZ-NH1	7.14	123.87	120.30
1	А	241	ARG	NE-CZ-NH2	-6.81	116.90	120.30
1	А	327	ARG	NE-CZ-NH2	-6.70	116.95	120.30
1	А	190	ASP	CB-CG-OD2	-6.34	112.59	118.30
1	А	136	ARG	NE-CZ-NH2	-6.06	117.27	120.30



Mol	Chain	Res	Type	Atoms	Ζ	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
1	А	190	ASP	CB-CG-OD1	5.92	123.63	118.30
1	А	200	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	А	78	MET	CA-CB-CG	5.57	122.76	113.30
1	А	131	ARG	NE-CZ-NH1	5.39	122.99	120.30
1	А	51	ASP	CB-CG-OD2	5.29	123.06	118.30
1	А	311	ASP	CB-CG-OD1	5.16	122.94	118.30
1	А	200	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	А	372	ARG	NE-CZ-NH2	-5.07	117.76	120.30

Continued from previous page...

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	61[A]	ARG	Mainchain

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	2807	0	2846	6	0
2	А	31	0	19	0	0
3	А	11	0	3	0	0
4	А	34	0	0	0	0
5	А	12	0	9	1	0
6	А	15	0	0	0	0
7	А	24	0	32	0	0
8	А	291	0	0	1	0
All	All	3225	0	2909	6	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 1.

All (6) 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:266[A]:GLU:OE2	8:A:503:HOH:O	1.64	1.14



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:315[B]:GLN:HE22	1:A:318:ARG:HE	1.48	0.60
1:A:315[B]:GLN:O	1:A:319[B]:GLU:HG3	2.02	0.60
1:A:315[B]:GLN:NE2	5:A:404:ACT:O	2.47	0.47
1:A:315[B]:GLN:NE2	1:A:315[B]:GLN:HA	2.35	0.41
1:A:203:GLY:N	1:A:204:PRO:CD	2.83	0.40

Continued from previous page...

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	Percentile	es
1	А	360/369~(98%)	348~(97%)	12 (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	А	298/296~(101%)	295~(99%)	3 (1%)	76 71	

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	51	ASP
	<i>a</i>	1	



Continued from previous page...

Mol	Chain	Res	Type	
1	А	246	ARG	
1	А	356	TYR	

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

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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)

13 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with |Z| > 2 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bo	ond leng	ths	B	ond ang	gles
	Type	Chain	nes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	SO4	A	410	-	$4,\!4,\!4$	0.37	0	6,6,6	0.25	0
7	GOL	А	413	-	$5,\!5,\!5$	0.40	0	$5,\!5,\!5$	0.40	0
5	ACT	А	404	-	3,3,3	1.07	0	3,3,3	0.45	0
6	SO4	А	406	-	4,4,4	0.66	0	6,6,6	0.66	0
6	SO4	А	411	-	4,4,4	0.41	0	6,6,6	0.33	0
7	GOL	А	409	-	$5,\!5,\!5$	0.54	0	$5,\!5,\!5$	0.54	0
4	A1AA2	А	403	-	$35,\!35,\!35$	1.20	4 (11%)	49,50,50	1.64	10 (20%)
3	ORO	А	402	-	9,11,11	1.14	1 (11%)	8,15,15	2.49	2 (25%)



Mol	Turne	Chain	Res	Link	Bo	ond leng	ths	Bond angles			
IVIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2	
5	ACT	А	407	-	$3,\!3,\!3$	0.93	0	$3,\!3,\!3$	0.49	0	
2	FMN	А	401	-	33,33,33	1.61	6 (18%)	$48,\!50,\!50$	1.55	10 (20%)	
7	GOL	А	408	-	$5,\!5,\!5$	0.46	0	$5,\!5,\!5$	0.52	0	
5	ACT	А	405	-	$3,\!3,\!3$	0.74	0	$3,\!3,\!3$	0.97	0	
7	GOL	А	412	-	$5,\!5,\!5$	0.31	0	$5,\!5,\!5$	0.69	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
7	GOL	А	413	-	-	0/4/4/4	-
7	GOL	А	409	-	-	2/4/4/4	-
4	A1AA2	А	403	-	-	1/31/31/31	0/2/2/2
3	ORO	А	402	-	-	4/4/4/4	0/1/1/1
2	FMN	А	401	-	-	2/18/18/18	0/3/3/3
7	GOL	А	408	-	-	2/4/4/4	-
7	GOL	А	412	-	-	0/4/4/4	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	Observed(Å)	Ideal(Å)
2	А	401	FMN	C9A-C5A	5.15	1.49	1.41
2	А	401	FMN	C4A-N5	3.19	1.37	1.30
2	А	401	FMN	C5A-N5	-2.55	1.34	1.39
2	А	401	FMN	C4-N3	-2.46	1.34	1.38
3	А	402	ORO	C2-N1	-2.34	1.33	1.38
2	А	401	FMN	C8-C7	2.29	1.46	1.40
4	А	403	A1AA2	O13-C14	-2.28	1.42	1.46
4	А	403	A1AA2	C28-N27	-2.28	1.39	1.43
2	А	401	FMN	C10-N10	2.25	1.42	1.37
4	А	403	A1AA2	F20-C17	2.22	1.39	1.33
4	А	403	A1AA2	C34-CL35	-2.10	1.68	1.73

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	А	402	ORO	C5-C4-N3	-5.48	117.69	124.08
4	А	403	A1AA2	C29-C28-N27	-4.18	116.89	122.12
2	А	401	FMN	O2-C2-N1	-4.08	115.06	121.83



Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	А	401	FMN	C4-C4A-N5	3.80	123.64	118.23
4	А	403	A1AA2	C31-C29-C28	-3.76	118.64	122.88
3	А	402	ORO	C6-C5-C4	3.62	119.06	116.73
4	А	403	A1AA2	N9-C7-N3	3.47	119.86	115.89
4	А	403	A1AA2	F20-C17-F19	3.12	113.83	106.87
2	А	401	FMN	C9A-C5A-N5	-2.78	119.41	122.43
4	А	403	A1AA2	C22-C23-C10	-2.74	120.89	123.50
2	А	401	FMN	O4-C4-C4A	-2.63	119.63	126.60
2	А	401	FMN	C5A-N5-C4A	2.50	122.23	118.07
4	А	403	A1AA2	C10-N9-C7	2.38	131.67	125.52
4	А	403	A1AA2	C12-O13-C14	2.30	123.97	119.52
2	А	401	FMN	O2-C2-N3	2.28	123.09	118.65
2	А	401	FMN	C10-N1-C2	2.26	121.43	116.90
2	А	401	FMN	C4A-C4-N3	2.20	118.77	113.19
4	А	403	A1AA2	C34-C28-C29	2.19	119.53	116.13
4	А	403	A1AA2	F30-C29-C28	2.14	119.54	117.70
2	А	401	FMN	C4A-C10-N1	-2.12	119.80	124.73
4	А	403	A1AA2	C11-C10-C23	2.05	119.63	117.75
2	А	401	FMN	O2P-P-O1P	2.02	118.59	110.68

Continued from previous page...

There are no chirality outliers.

Mol	Chain	Res	Type	Atoms
3	А	402	ORO	N1-C6-C7-O72
3	А	402	ORO	C5-C6-C7-O71
3	А	402	ORO	C5-C6-C7-O72
7	А	408	GOL	O1-C1-C2-C3
7	А	409	GOL	O1-C1-C2-C3
3	А	402	ORO	N1-C6-C7-O71
7	А	408	GOL	O1-C1-C2-O2
7	А	409	GOL	O1-C1-C2-O2
2	А	401	FMN	C4'-C5'-O5'-P
4	А	403	A1AA2	C16-C14-O13-C12
2	А	401	FMN	C5'-O5'-P-O1P

All (11) torsion outliers are listed below:

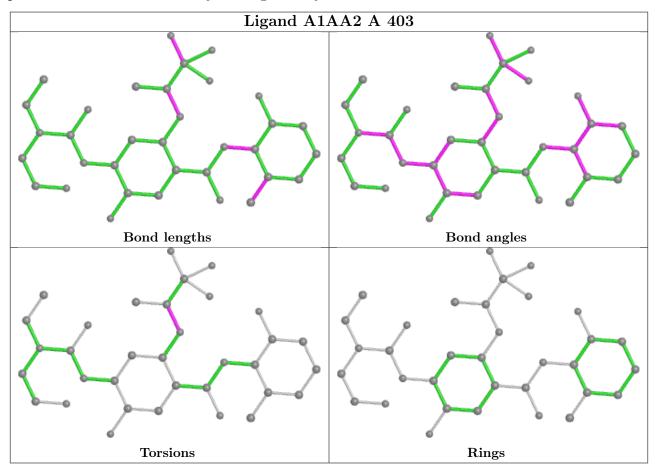
There are no ring outliers.

1 monomer is involved in 1 short contact:

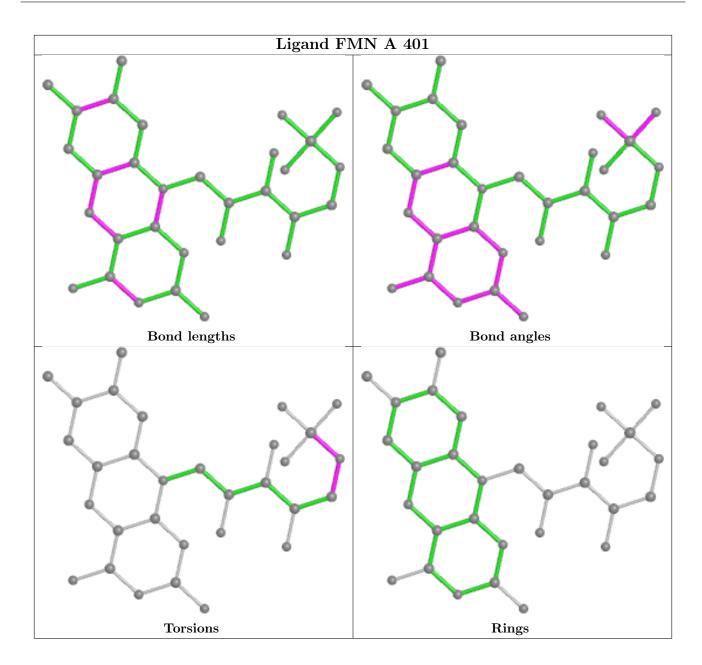
M	bl	Chain	Res	Type	Clashes	Symm-Clashes
5		А	404	ACT	1	0



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 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.







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^{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	А	355/369~(96%)	0.15	23 (6%) 18 25	10, 17, 38, 77	7 (1%)

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	216	PRO	6.2
1	А	70	ARG	5.7
1	А	189	VAL	4.9
1	А	225	GLN	4.6
1	А	73	PHE	4.5
1	А	31	ALA	4.0
1	А	69	PRO	3.8
1	А	228	ALA	3.8
1	А	33	GLY	3.7
1	А	37	PHE	3.6
1	А	226	GLY	3.5
1	А	74	GLN	3.4
1	А	215	SER	2.8
1	А	214	SER	2.7
1	А	94	ILE	2.6
1	А	187	THR	2.6
1	А	49	LEU	2.4
1	А	32	THR	2.3
1	А	355	LEU	2.2
1	А	84	LEU	2.1
1	А	353	VAL	2.1
1	А	95	ALA	2.1
1	А	92	VAL	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.

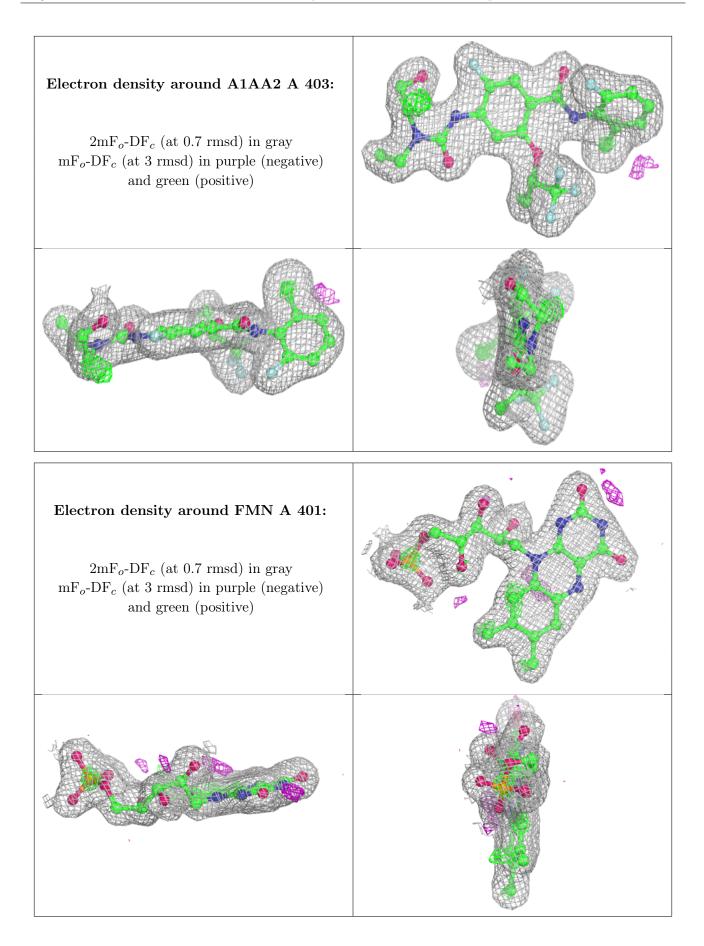
6.4 Ligands (i)

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, 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
7	GOL	А	409	6/6	0.69	0.31	54,63,80,90	0
7	GOL	A	412	6/6	0.71	0.33	59,75,84,96	0
6	SO4	A	411	5/5	0.75	0.19	$65,\!67,\!93,\!102$	0
6	SO4	A	406	5/5	0.84	0.15	51,60,73,140	0
7	GOL	A	413	6/6	0.84	0.15	62,66,81,82	0
7	GOL	A	408	6/6	0.88	0.17	48,54,57,61	0
5	ACT	A	407	4/4	0.89	0.14	43,44,46,46	0
5	ACT	A	404	4/4	0.92	0.29	40,40,47,53	0
4	A1AA2	A	403	34/34	0.92	0.12	21,29,34,36	0
6	SO4	A	410	5/5	0.93	0.16	53,63,72,80	0
3	ORO	А	402	11/11	0.96	0.12	23,26,29,30	0
2	FMN	A	401	31/31	0.97	0.17	16,22,26,26	0
5	ACT	А	405	4/4	0.97	0.15	32,40,40,65	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.







6.5 Other polymers (i)

There are no such residues in this entry.

