



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:18 PM UTC

PDB ID : 4P0W / pdb_00004p0w
Title : Human farnesyl diphosphate synthase in complex with Arenarone and zole-dronate
Authors : Liu, Y.L.; Oldfield, E.
Deposited on : 2014-02-23
Resolution : 2.41 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

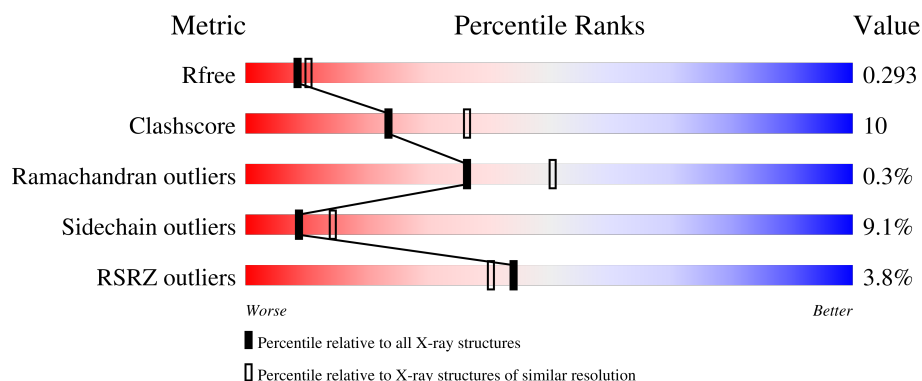
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

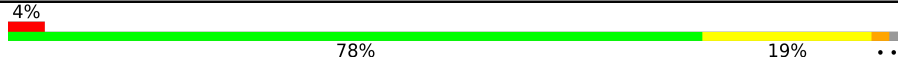
The reported resolution of this entry is 2.41 Å.

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}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	348	

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
2	1XH	A	500	-	-	X	-

2 Entry composition [i](#)

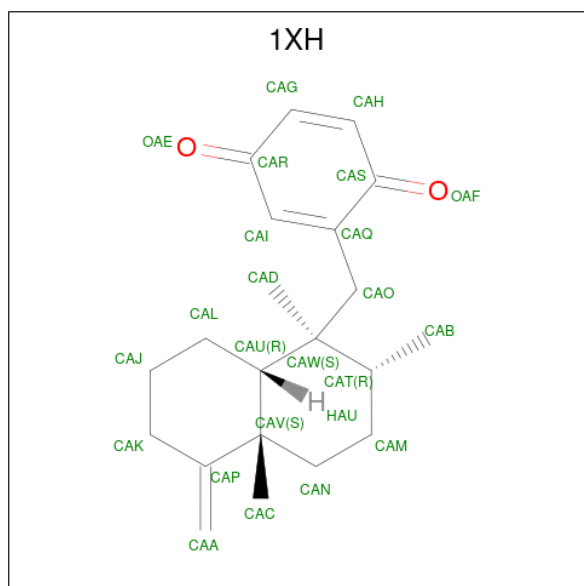
There are 5 unique types of molecules in this entry. The entry contains 2904 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Farnesyl pyrophosphate synthase.

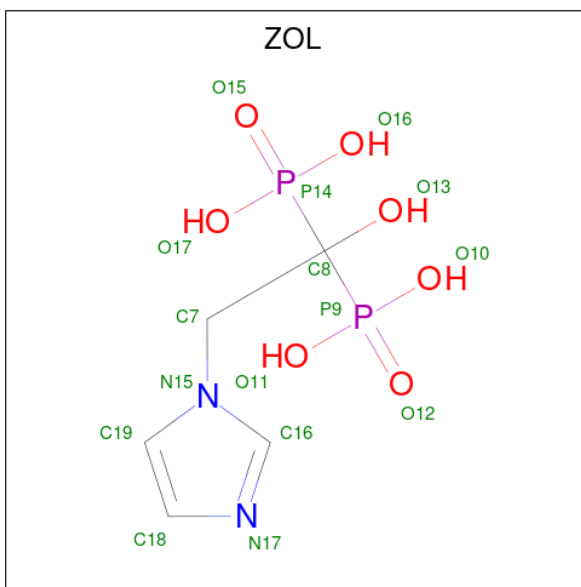
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	346	2806	1806	464	524	12	0	0	0

- Molecule 2 is 2-{{[(1S,2R,4aS,8aR)-1,2,4a-trimethyl-5-methylenedecahydronaphthalen-1-yl]methyl}cyclohexa-2,5-diene-1,4-dione (CCD ID: 1XH) (formula: C₂₁H₂₈O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	23	21	2	0	0

- Molecule 3 is ZOLEDRONIC ACID (CCD ID: ZOL) (formula: C₅H₁₀N₂O₇P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			16	5	2	7	2		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Mg	0	0
			3	3		

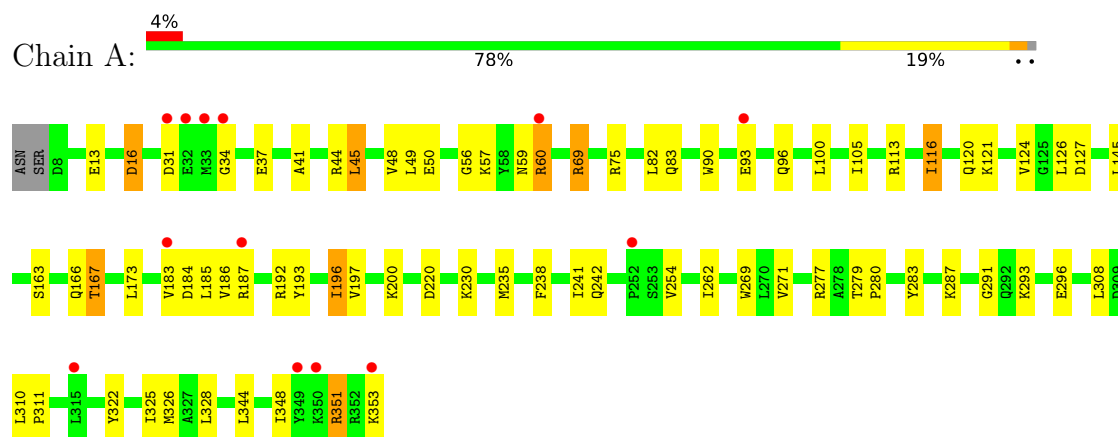
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	56	Total	O	0	0
			56	56		

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: Farnesyl pyrophosphate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	111.24Å 111.24Å 67.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	57.58 – 2.41 57.58 – 2.41	Depositor EDS
% Data completeness (in resolution range)	97.2 (57.58-2.41) 97.3 (57.58-2.41)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	27.23 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0069	Depositor
R, R_{free}	0.204 , 0.289 0.208 , 0.293	Depositor DCC
R_{free} test set	835 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	42.1	Xtriage
Anisotropy	0.224	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2904	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.85% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 1XH, ZOL, MG

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	1.02	3/2865 (0.1%)	1.08	3/3876 (0.1%)

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	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	41	ALA	CA-C	-6.00	1.45	1.52
1	A	105	ILE	C-O	-5.38	1.18	1.24
1	A	196	ILE	N-CA	-5.18	1.40	1.46

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	325	ILE	CB-CA-C	-5.26	105.04	112.14
1	A	279	THR	CA-C-N	-5.02	114.44	119.56
1	A	279	THR	C-N-CA	-5.02	114.44	119.56

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	183	VAL	Peptide

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Mol	Chain	Res	Type	Group
1	A	184	ASP	Peptide
1	A	34	GLY	Peptide

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	2806	0	2791	47	0
2	A	23	0	27	32	0
3	A	16	0	6	0	0
4	A	3	0	0	0	0
5	A	56	0	0	5	0
All	All	2904	0	2824	58	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 10.

All (58) 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:96:GLN:HE21	2:A:500:1XH:HAU	1.16	1.09
2:A:500:1XH:HAAA	5:A:652:HOH:O	1.54	1.06
1:A:93:GLU:CG	2:A:500:1XH:OAE	2.15	0.95
1:A:56:GLY:HA3	2:A:500:1XH:HAB	1.56	0.86
1:A:93:GLU:HB2	2:A:500:1XH:OAE	1.75	0.85
1:A:93:GLU:CB	2:A:500:1XH:OAE	2.26	0.84
1:A:60:ARG:H	2:A:500:1XH:HAH	1.51	0.76
1:A:96:GLN:NE2	2:A:500:1XH:HAU	1.97	0.75
1:A:96:GLN:HE21	2:A:500:1XH:CAU	1.97	0.73
2:A:500:1XH:OAF	2:A:500:1XH:HAL	1.88	0.73
1:A:163:SER:O	1:A:167:THR:HG23	1.88	0.72
2:A:500:1XH:HAD	2:A:500:1XH:HANA	1.71	0.71
2:A:500:1XH:OAF	2:A:500:1XH:CAL	2.41	0.68
1:A:113:ARG:HH21	2:A:500:1XH:HAO	1.60	0.66
1:A:93:GLU:HG2	2:A:500:1XH:OAE	1.96	0.65
1:A:113:ARG:NH2	2:A:500:1XH:HAO	2.12	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:500:1XH:CAA	5:A:652:HOH:O	2.25	0.63
1:A:57:LYS:H	2:A:500:1XH:HADA	1.64	0.63
1:A:56:GLY:CA	2:A:500:1XH:HAB	2.28	0.62
1:A:60:ARG:N	2:A:500:1XH:HAH	2.16	0.60
1:A:93:GLU:HG3	2:A:500:1XH:OAE	2.00	0.59
1:A:96:GLN:NE2	2:A:500:1XH:HAC	2.17	0.59
1:A:235:MET:HE1	1:A:344:LEU:HG	1.85	0.58
1:A:163:SER:O	1:A:167:THR:CG2	2.50	0.57
1:A:163:SER:HA	1:A:166:GLN:HE21	1.68	0.57
1:A:59:ASN:HB2	1:A:60:ARG:CZ	2.37	0.55
2:A:500:1XH:HAD	2:A:500:1XH:CAN	2.32	0.54
1:A:116:ILE:HG22	1:A:120:GLN:HB2	1.89	0.54
1:A:242:GLN:NE2	1:A:351:ARG:HE	2.05	0.53
2:A:500:1XH:CAT	2:A:500:1XH:HACA	2.39	0.52
1:A:93:GLU:CD	1:A:93:GLU:C	2.78	0.51
1:A:310:LEU:N	1:A:311:PRO:CD	2.75	0.50
1:A:60:ARG:CB	2:A:500:1XH:HAH	2.42	0.50
1:A:13:GLU:OE2	1:A:69:ARG:HD2	2.12	0.49
1:A:96:GLN:NE2	2:A:500:1XH:CAU	2.68	0.49
2:A:500:1XH:HAN	5:A:652:HOH:O	2.12	0.49
1:A:254:VAL:HG12	1:A:353:LYS:HB3	1.94	0.48
1:A:238:PHE:CD2	1:A:348:ILE:HD12	2.49	0.47
1:A:60:ARG:HB2	2:A:500:1XH:HAH	1.96	0.47
1:A:60:ARG:H	2:A:500:1XH:CAH	2.25	0.47
1:A:192:ARG:O	1:A:196:ILE:HG13	2.15	0.47
2:A:500:1XH:HACB	5:A:652:HOH:O	2.15	0.46
1:A:322:TYR:CE1	1:A:326:MET:HE3	2.50	0.46
1:A:280:PRO:O	1:A:283:TYR:HB3	2.15	0.46
2:A:500:1XH:HAAA	2:A:500:1XH:HACB	1.58	0.46
1:A:100:LEU:HD11	2:A:500:1XH:HAC	1.98	0.46
1:A:90:TRP:CZ3	1:A:93:GLU:HG3	2.52	0.45
1:A:16:ASP:OD1	1:A:16:ASP:N	2.50	0.44
1:A:124:VAL:O	1:A:127:ASP:HB2	2.17	0.44
1:A:262:ILE:O	1:A:291:GLY:HA2	2.17	0.44
1:A:193:TYR:CZ	1:A:197:VAL:HG11	2.52	0.44
1:A:241:ILE:HG12	1:A:269:TRP:CD1	2.53	0.44
1:A:196:ILE:O	1:A:200:LYS:HG2	2.19	0.43
1:A:220:ASP:OD1	1:A:220:ASP:N	2.52	0.42
1:A:167:THR:HG22	5:A:631:HOH:O	2.18	0.42
1:A:45:LEU:HD22	1:A:49:LEU:CD1	2.51	0.41
2:A:500:1XH:HADA	2:A:500:1XH:HABA	1.71	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:ARG:O	1:A:48:VAL:HG23	2.20	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	344/348 (99%)	330 (96%)	13 (4%)	1 (0%)	36 50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	185	LEU

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	297/299 (99%)	270 (91%)	27 (9%)	9 14

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	ASP
1	A	31	ASP

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Mol	Chain	Res	Type
1	A	37	GLU
1	A	45	LEU
1	A	50	GLU
1	A	60	ARG
1	A	69	ARG
1	A	75	ARG
1	A	82	LEU
1	A	83	GLN
1	A	116	ILE
1	A	121	LYS
1	A	126	LEU
1	A	145	LEU
1	A	167	THR
1	A	173	LEU
1	A	186	VAL
1	A	187	ARG
1	A	230	LYS
1	A	271	VAL
1	A	277	ARG
1	A	287	LYS
1	A	293	LYS
1	A	296	GLU
1	A	308	LEU
1	A	328	LEU
1	A	351	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	96	GLN
1	A	162	GLN
1	A	166	GLN
1	A	242	GLN
1	A	289	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 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	$\# Z > 2$
3	ZOL	A	501	4	16,16,16	4.01	9 (56%)	23,26,26	2.30	7 (30%)
2	1XH	A	500	-	25,25,25	5.62	18 (72%)	36,39,39	4.54	25 (69%)

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	ZOL	A	501	4	-	2/23/23/23	0/1/1/1
2	1XH	A	500	-	-	2/5/54/54	0/3/3/3

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	1XH	CAW-CAT	-14.69	1.31	1.55
3	A	501	ZOL	P9-O12	11.47	1.69	1.50
2	A	500	1XH	CAN-CAV	-9.45	1.36	1.54
2	A	500	1XH	CAW-CAU	9.09	1.70	1.56
2	A	500	1XH	CAV-CAU	8.32	1.68	1.56
2	A	500	1XH	CAA-CAP	7.15	1.44	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	1XH	CAO-CAW	6.61	1.69	1.56
3	A	501	ZOL	P9-C8	-6.49	1.80	1.85
2	A	500	1XH	CAO-CAQ	-6.38	1.44	1.50
2	A	500	1XH	CAB-CAT	-6.25	1.39	1.53
2	A	500	1XH	CAM-CAT	-6.20	1.44	1.53
3	A	501	ZOL	P14-C8	-5.85	1.81	1.85
2	A	500	1XH	CAV-CAP	-5.30	1.39	1.53
2	A	500	1XH	CAL-CAU	4.23	1.60	1.53
2	A	500	1XH	CAI-CAQ	3.99	1.44	1.34
2	A	500	1XH	CAG-CAR	-3.79	1.36	1.46
2	A	500	1XH	CAQ-CAS	-3.46	1.44	1.49
3	A	501	ZOL	P14-O15	-3.38	1.44	1.50
3	A	501	ZOL	P14-O17	-2.71	1.47	1.54
3	A	501	ZOL	P9-O10	-2.68	1.47	1.54
2	A	500	1XH	CAK-CAP	-2.64	1.46	1.50
2	A	500	1XH	CAN-CAM	-2.55	1.48	1.53
3	A	501	ZOL	C7-C8	-2.51	1.50	1.54
2	A	500	1XH	CAJ-CAK	2.28	1.59	1.52
3	A	501	ZOL	O13-C8	-2.27	1.41	1.44
3	A	501	ZOL	P9-O11	-2.19	1.49	1.54
2	A	500	1XH	CAC-CAV	2.04	1.57	1.54

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	1XH	CAB-CAT-CAM	-10.14	95.54	110.33
2	A	500	1XH	CAI-CAQ-CAS	-8.50	112.60	118.66
2	A	500	1XH	CAU-CAV-CAP	8.40	122.93	107.39
2	A	500	1XH	CAG-CAR-CAI	8.11	125.34	117.12
2	A	500	1XH	OAE-CAR-CAG	-7.85	108.92	121.56
2	A	500	1XH	CAN-CAV-CAU	-7.54	100.02	108.36
2	A	500	1XH	CAU-CAW-CAT	-5.48	100.81	108.73
2	A	500	1XH	CAC-CAV-CAN	5.09	117.18	109.43
2	A	500	1XH	CAN-CAM-CAT	5.05	118.89	112.38
2	A	500	1XH	CAO-CAW-CAU	5.04	119.85	109.80
2	A	500	1XH	CAM-CAN-CAV	-4.94	102.27	112.78
3	A	501	ZOL	O11-P9-O10	4.91	122.08	108.24
3	A	501	ZOL	C7-N15-C19	-4.72	118.27	127.02
2	A	500	1XH	CAO-CAQ-CAS	4.69	126.77	118.56
2	A	500	1XH	CAH-CAG-CAR	-4.52	115.11	121.25
2	A	500	1XH	CAD-CAW-CAT	-4.48	104.98	110.83
3	A	501	ZOL	C7-N15-C16	4.18	133.44	125.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	ZOL	O10-P9-O12	-4.11	102.88	112.96
2	A	500	1XH	OAF-CAS-CAH	-4.08	114.25	121.58
2	A	500	1XH	CAV-CAP-CAA	-3.80	116.97	123.98
2	A	500	1XH	CAN-CAV-CAP	-3.55	103.12	111.19
2	A	500	1XH	CAL-CAU-CAW	3.52	119.03	114.31
3	A	501	ZOL	O15-P14-C8	-3.38	101.15	109.82
2	A	500	1XH	CAH-CAS-CAQ	3.37	123.43	117.48
2	A	500	1XH	OAE-CAR-CAI	2.97	125.65	121.43
2	A	500	1XH	CAK-CAP-CAV	2.77	118.58	115.19
2	A	500	1XH	CAB-CAT-CAW	2.69	117.41	114.38
3	A	501	ZOL	C8-C7-N15	-2.57	110.41	113.67
2	A	500	1XH	CAW-CAU-CAV	-2.41	112.89	115.86
3	A	501	ZOL	O17-P14-O15	2.22	118.41	112.96
2	A	500	1XH	CAJ-CAK-CAP	2.11	118.50	111.68
2	A	500	1XH	CAK-CAP-CAA	2.06	124.42	121.32

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	500	1XH	CAW-CAO-CAQ-CAI
2	A	500	1XH	CAW-CAO-CAQ-CAS
3	A	501	ZOL	C8-C7-N15-C16
3	A	501	ZOL	C8-C7-N15-C19

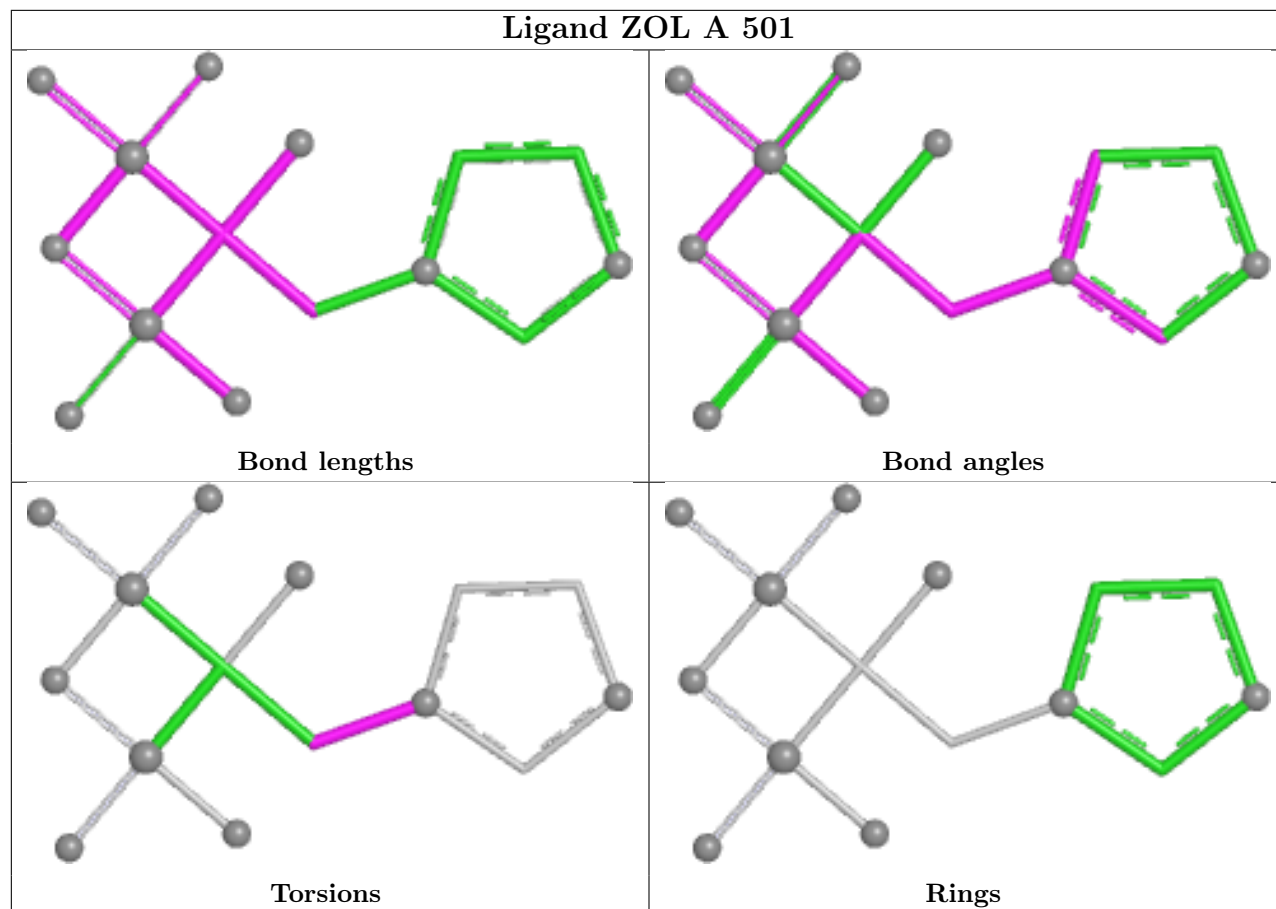
There are no ring outliers.

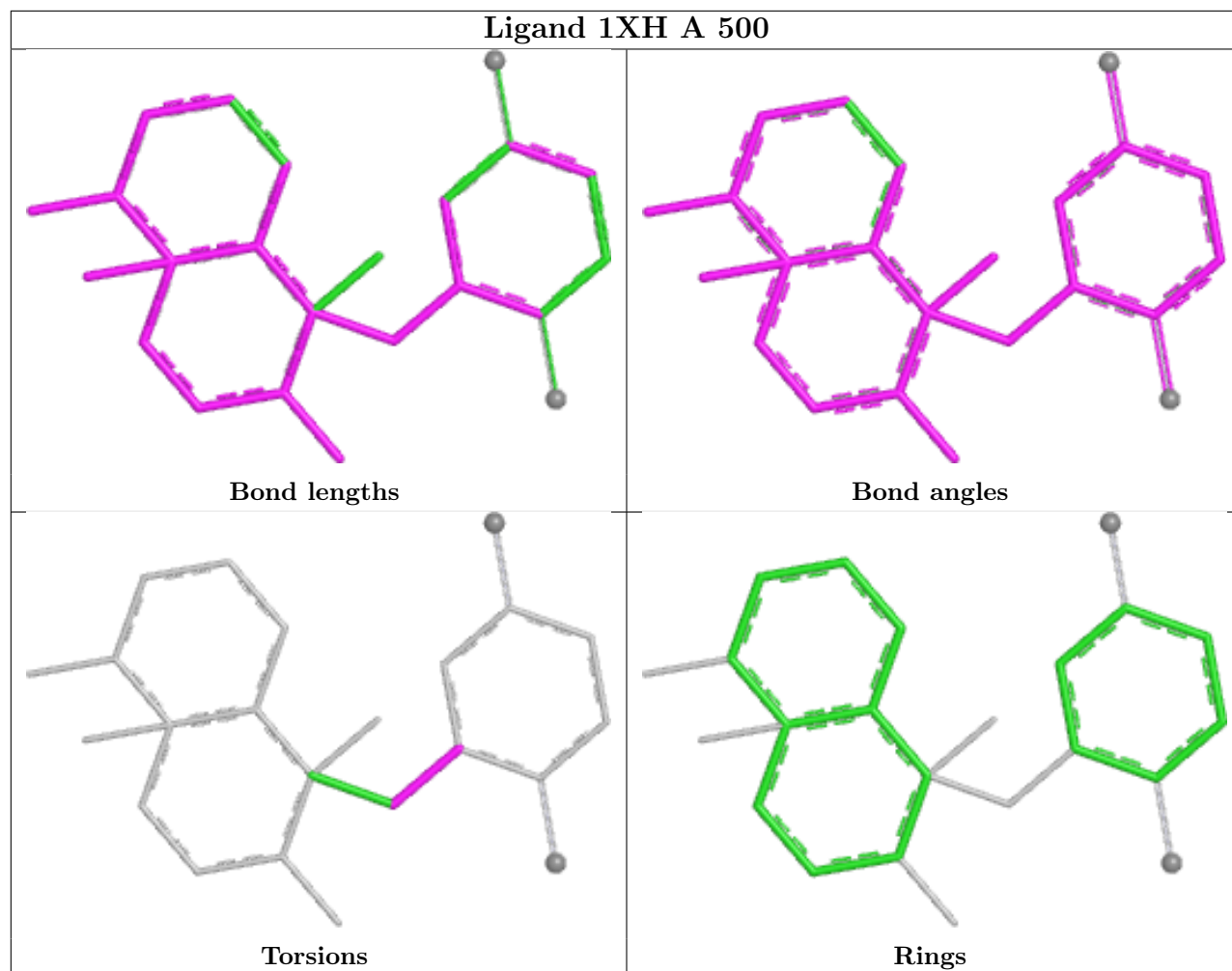
1 monomer is involved in 32 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	1XH	32	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.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	346/348 (99%)	0.17	13 (3%) 44 40	27, 47, 78, 113	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	60	ARG	4.5
1	A	93	GLU	4.3
1	A	349	TYR	3.8
1	A	31	ASP	2.7
1	A	183	VAL	2.6
1	A	34	GLY	2.6
1	A	350	LYS	2.4
1	A	353	LYS	2.2
1	A	315	LEU	2.2
1	A	187	ARG	2.1
1	A	32	GLU	2.1
1	A	252	PRO	2.0
1	A	33	MET	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](#)

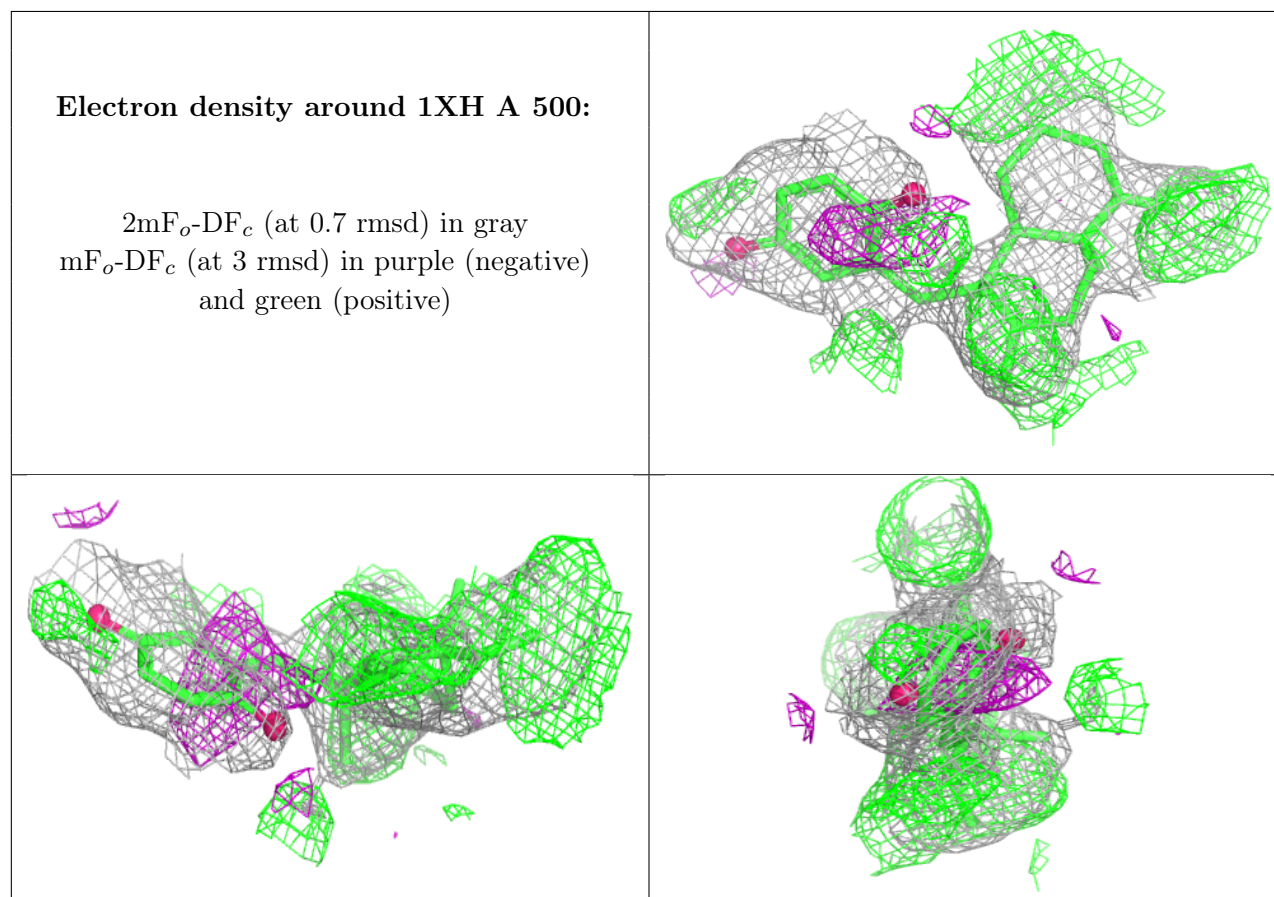
There are no oligosaccharides in this entry.

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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

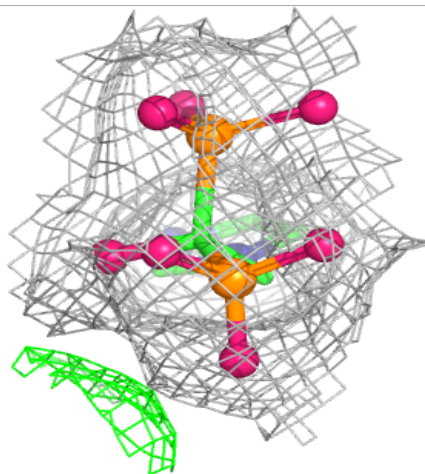
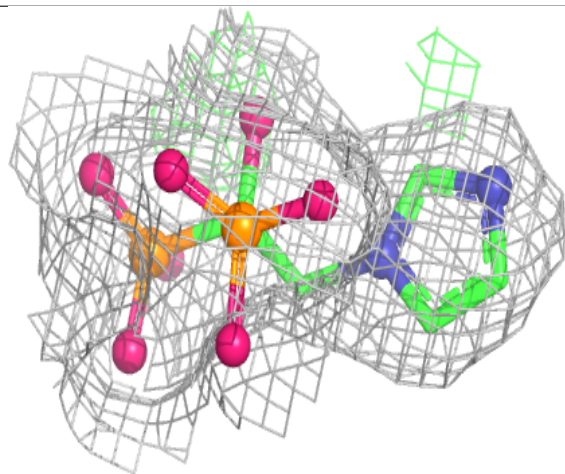
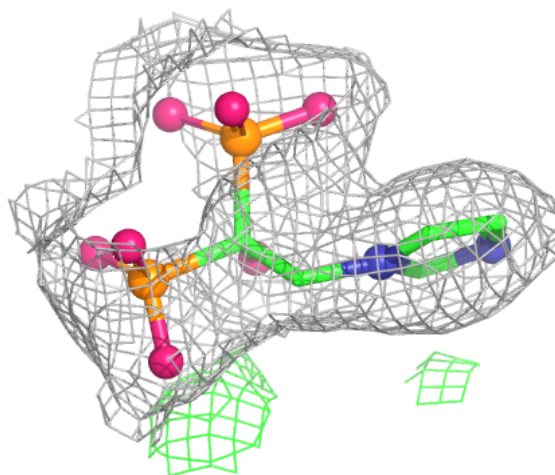
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	1XH	A	500	23/23	0.67	0.24	27,46,52,63	0
3	ZOL	A	501	16/16	0.98	0.05	27,33,34,35	0
4	MG	A	502	1/1	0.99	0.03	32,32,32,32	0
4	MG	A	503	1/1	0.99	0.02	34,34,34,34	0
4	MG	A	504	1/1	1.00	0.04	14,14,14,14	0

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 ZOL A 501:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.