



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 2, 2024 – 02:34 pm GMT

PDB ID : 4UMX
Title : IDH1 R132H in complex with cpd 1
Authors : Mathieu, M.; Marquette, J.P.
Deposited on : 2014-05-22
Resolution : 1.88 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

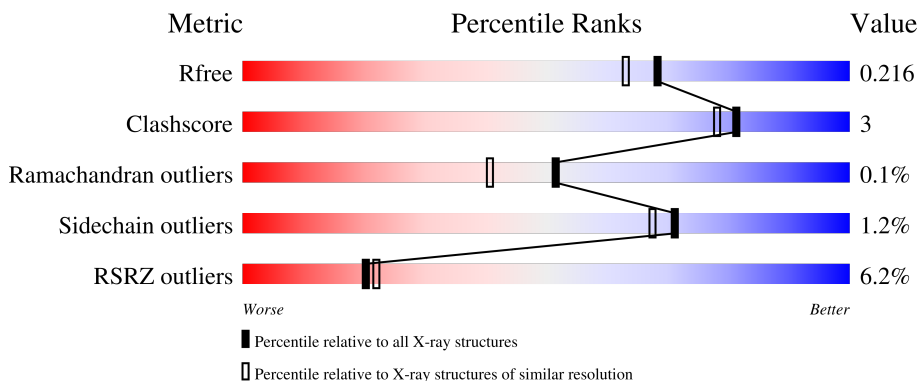
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.88 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	9470 (1.90-1.86)
Clashscore	141614	10282 (1.90-1.86)
Ramachandran outliers	138981	10152 (1.90-1.86)
Sidechain outliers	138945	10152 (1.90-1.86)
RSRZ outliers	127900	9303 (1.90-1.86)

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	425	 9% 92% 5%
1	B	425	 3% 88% 7% 5%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7374 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.

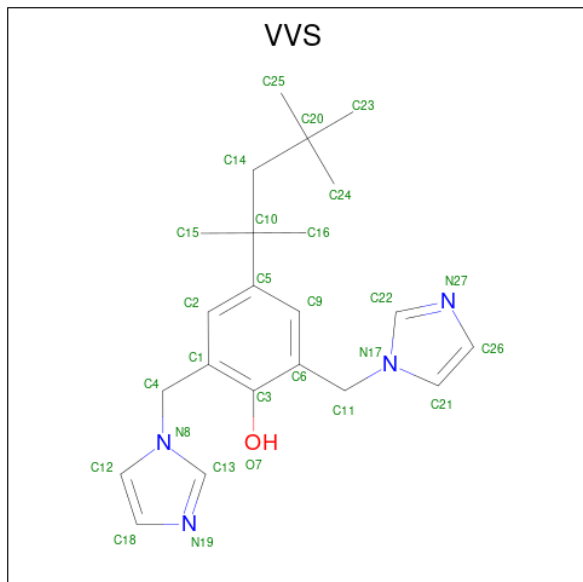
- Molecule 1 is a protein called ISOCITRATE DEHYDROGENASE [NADP] CYTOPLASMIC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	412	3265	2077	553	617	18	0	0	0
1	B	405	3214	2046	545	605	18	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

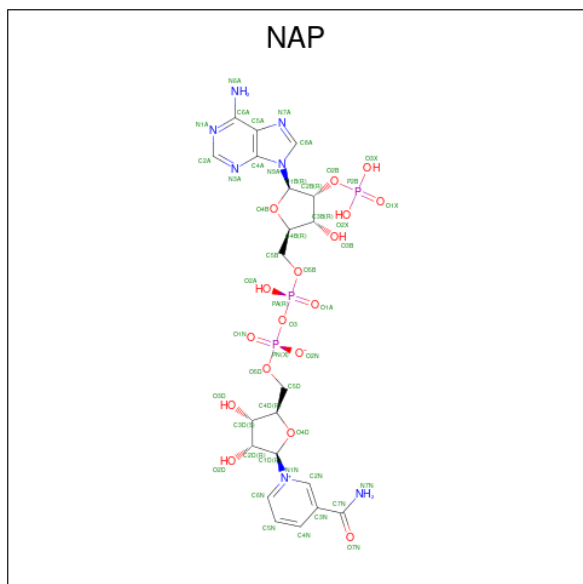
Chain	Residue	Modelled	Actual	Comment	Reference
A	415	SER	-	expression tag	UNP O75874
A	416	LEU	-	expression tag	UNP O75874
A	417	GLU	-	expression tag	UNP O75874
A	418	HIS	-	expression tag	UNP O75874
A	419	HIS	-	expression tag	UNP O75874
A	420	HIS	-	expression tag	UNP O75874
A	421	HIS	-	expression tag	UNP O75874
A	422	HIS	-	expression tag	UNP O75874
A	423	HIS	-	expression tag	UNP O75874
A	424	HIS	-	expression tag	UNP O75874
A	425	HIS	-	expression tag	UNP O75874
A	132	HIS	ARG	engineered mutation	UNP O75874
B	415	SER	-	expression tag	UNP O75874
B	416	LEU	-	expression tag	UNP O75874
B	417	GLU	-	expression tag	UNP O75874
B	418	HIS	-	expression tag	UNP O75874
B	419	HIS	-	expression tag	UNP O75874
B	420	HIS	-	expression tag	UNP O75874
B	421	HIS	-	expression tag	UNP O75874
B	422	HIS	-	expression tag	UNP O75874
B	423	HIS	-	expression tag	UNP O75874
B	424	HIS	-	expression tag	UNP O75874
B	425	HIS	-	expression tag	UNP O75874
B	132	HIS	ARG	engineered mutation	UNP O75874

- Molecule 2 is 2,6-bis(1H-imidazol-1-ylmethyl)-4-(2,4,4-trimethylpentan-2-yl)phenol (three-letter code: VVS) (formula: C₂₂H₃₀N₄O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	27	22	4	1	0	0

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



Continued from previous page...

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	N	O			P
3	B	1	48	21	7	17	3	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	B	1	6	3	3	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	377	377	377	0	0
5	B	389	389	389	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	81.41Å 81.41Å 305.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	78.67 – 1.88 78.67 – 1.88	Depositor EDS
% Data completeness (in resolution range)	100.0 (78.67-1.88) 100.0 (78.67-1.88)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.76 (at 1.87Å)	Xtrriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.182 , 0.211 0.184 , 0.216	Depositor DCC
R_{free} test set	4278 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	25.6	Xtrriage
Anisotropy	0.097	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7374	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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: NAP, GOL, VVS

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	0.53	0/3334	0.63	0/4496
1	B	0.55	0/3281	0.62	0/4422
All	All	0.54	0/6615	0.63	0/8918

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3265	0	3233	11	0
1	B	3214	0	3189	24	0
2	A	27	0	30	0	0
3	A	48	0	25	1	0
3	B	48	0	25	1	0
4	B	6	0	8	2	0
5	A	377	0	0	0	0
5	B	389	0	0	4	0
All	All	7374	0	6510	34	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 3.

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:B:135:TYR:CB	1:B:139:TYR:HD2	2.04	0.71
1:A:208:TYR:HB3	1:A:254:MET:CE	2.24	0.67
1:B:315:HIS:HA	1:B:318:MET:HE2	1.79	0.65
1:B:135:TYR:HB2	1:B:139:TYR:HD2	1.62	0.64
1:A:315:HIS:HA	1:A:318:MET:HE2	1.80	0.64
1:B:94:SER:OG	4:B:1416:GOL:H31	1.99	0.62
1:A:208:TYR:HB3	1:A:254:MET:HE1	1.85	0.57
1:B:135:TYR:HB3	1:B:139:TYR:CD2	2.41	0.56
1:A:121:VAL:HB	1:A:124:TRP:CE2	2.41	0.56
1:B:135:TYR:HB3	1:B:139:TYR:HD2	1.70	0.55
1:B:182:MET:SD	1:B:273:ASP:HB3	2.47	0.55
1:A:75:THR:O	3:A:1416:NAP:H2N	2.09	0.53
1:B:135:TYR:CB	1:B:139:TYR:CD2	2.89	0.53
1:B:255:VAL:O	1:B:259:MET:HG2	2.08	0.52
1:B:96:ASN:HB2	4:B:1416:GOL:H11	1.91	0.52
1:B:141:ALA:HB1	1:B:180:MET:HE2	1.92	0.52
1:B:210:SER:HA	1:B:249:ARG:O	2.11	0.51
1:B:330:ILE:HD12	1:B:363:SER:HB3	1.91	0.51
1:B:75:THR:O	3:B:1415:NAP:H2N	2.11	0.50
1:B:141:ALA:HB1	1:B:180:MET:CE	2.42	0.50
1:B:286:GLY:HA2	1:B:287:SER:C	2.33	0.49
1:B:292:THR:HB	5:B:2278:HOH:O	2.14	0.48
1:A:330:ILE:HD12	1:A:363:SER:HB3	1.95	0.47
1:B:22:ILE:HD11	1:B:327:THR:HB	1.97	0.47
1:A:210:SER:HG	1:A:267:TRP:HE1	1.63	0.46
1:A:315:HIS:HA	1:A:318:MET:CE	2.43	0.46
1:A:156:TYR:CE2	1:B:147:PRO:HD2	2.51	0.46
1:B:315:HIS:HA	1:B:318:MET:CE	2.46	0.45
1:B:338:ARG:NH1	5:B:2270:HOH:O	2.50	0.44
1:A:208:TYR:HB3	1:A:254:MET:HE3	1.98	0.44
1:B:338:ARG:HG2	5:B:2270:HOH:O	2.18	0.43
1:B:133:HIS:HB2	1:B:192:PHE:CE2	2.55	0.41
1:A:255:VAL:O	1:A:259:MET:HG2	2.21	0.40
1:B:294:VAL:HG23	5:B:2270:HOH:O	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	410/425 (96%)	401 (98%)	8 (2%)	1 (0%)	47	37
1	B	401/425 (94%)	393 (98%)	8 (2%)	0	100	100
All	All	811/850 (95%)	794 (98%)	16 (2%)	1 (0%)	51	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	285	TYR

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	348/361 (96%)	342 (98%)	6 (2%)	60	54
1	B	343/361 (95%)	341 (99%)	2 (1%)	86	86
All	All	691/722 (96%)	683 (99%)	8 (1%)	71	67

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	171	ASN
1	A	251	ILE
1	A	260	LYS
1	A	270	LYS
1	A	312	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	385	ASN
1	B	171	ASN
1	B	312	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	348	ASN
1	B	357	ASN

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](#)

4 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	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	B	1416	-	5,5,5	0.06	0	5,5,5	0.38	0
3	NAP	A	1416	-	45,52,52	0.96	2 (4%)	56,80,80	0.83	2 (3%)
3	NAP	B	1415	-	45,52,52	1.10	2 (4%)	56,80,80	0.83	3 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	VVS	A	1415	-	27,29,29	0.54	0	39,43,43	0.83	1 (2%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	1416	-	-	2/4/4/4	-
3	NAP	A	1416	-	-	8/31/67/67	0/5/5/5
3	NAP	B	1415	-	-	9/31/67/67	0/5/5/5
2	VVS	A	1415	-	-	1/20/20/20	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1415	NAP	C2N-N1N	5.40	1.41	1.35
3	A	1416	NAP	C2N-N1N	3.76	1.39	1.35
3	A	1416	NAP	O7N-C7N	2.80	1.29	1.24
3	B	1415	NAP	O7N-C7N	2.79	1.29	1.24

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1415	NAP	O4D-C1D-C2D	-2.76	102.90	106.93
3	A	1416	NAP	C2N-N1N-C1D	-2.62	113.31	119.14
3	B	1415	NAP	C6N-N1N-C2N	-2.15	120.01	121.97
2	A	1415	VVS	C20-C14-C10	2.15	131.64	123.91
3	A	1416	NAP	O4D-C1D-C2D	-2.07	103.90	106.93
3	B	1415	NAP	C5A-C6A-N6A	2.01	123.41	120.35

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1416	NAP	C5D-O5D-PN-O2N
3	A	1416	NAP	O4D-C1D-N1N-C2N
3	A	1416	NAP	O4D-C1D-N1N-C6N
3	A	1416	NAP	C2D-C1D-N1N-C2N
3	B	1415	NAP	C5B-O5B-PA-O1A

Continued on next page...

Continued from previous page...

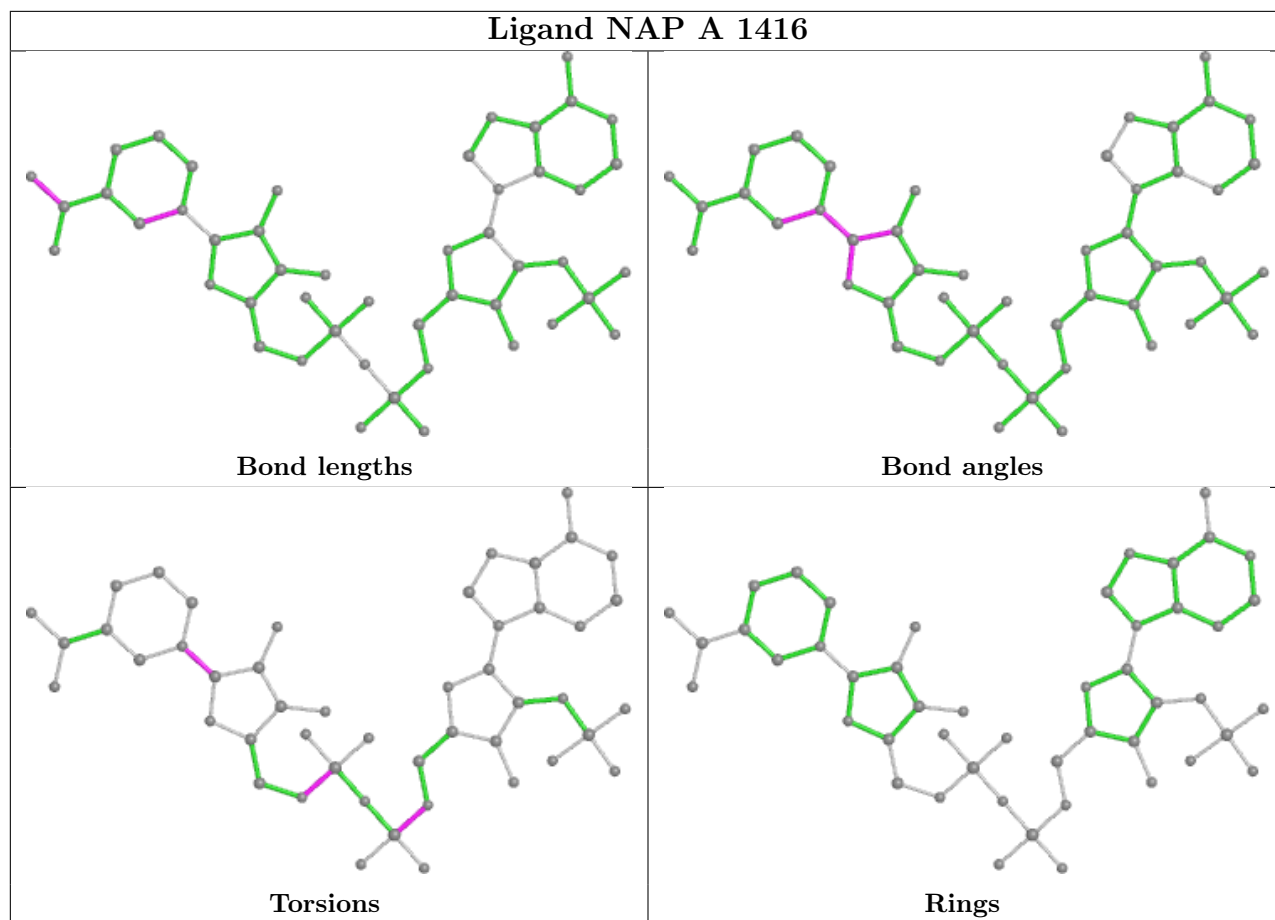
Mol	Chain	Res	Type	Atoms
3	B	1415	NAP	C5D-O5D-PN-O2N
3	B	1415	NAP	O4D-C1D-N1N-C2N
3	B	1415	NAP	O4D-C1D-N1N-C6N
3	B	1415	NAP	C2D-C1D-N1N-C2N
4	B	1416	GOL	C1-C2-C3-O3
4	B	1416	GOL	O2-C2-C3-O3
2	A	1415	VVS	N17-C11-C6-C3
3	A	1416	NAP	C5D-O5D-PN-O3
3	B	1415	NAP	C5D-O5D-PN-O3
3	A	1416	NAP	C5D-O5D-PN-O1N
3	B	1415	NAP	C5D-O5D-PN-O1N
3	A	1416	NAP	C2D-C1D-N1N-C6N
3	B	1415	NAP	C5B-O5B-PA-O3
3	B	1415	NAP	C2D-C1D-N1N-C6N
3	A	1416	NAP	C5B-O5B-PA-O1A

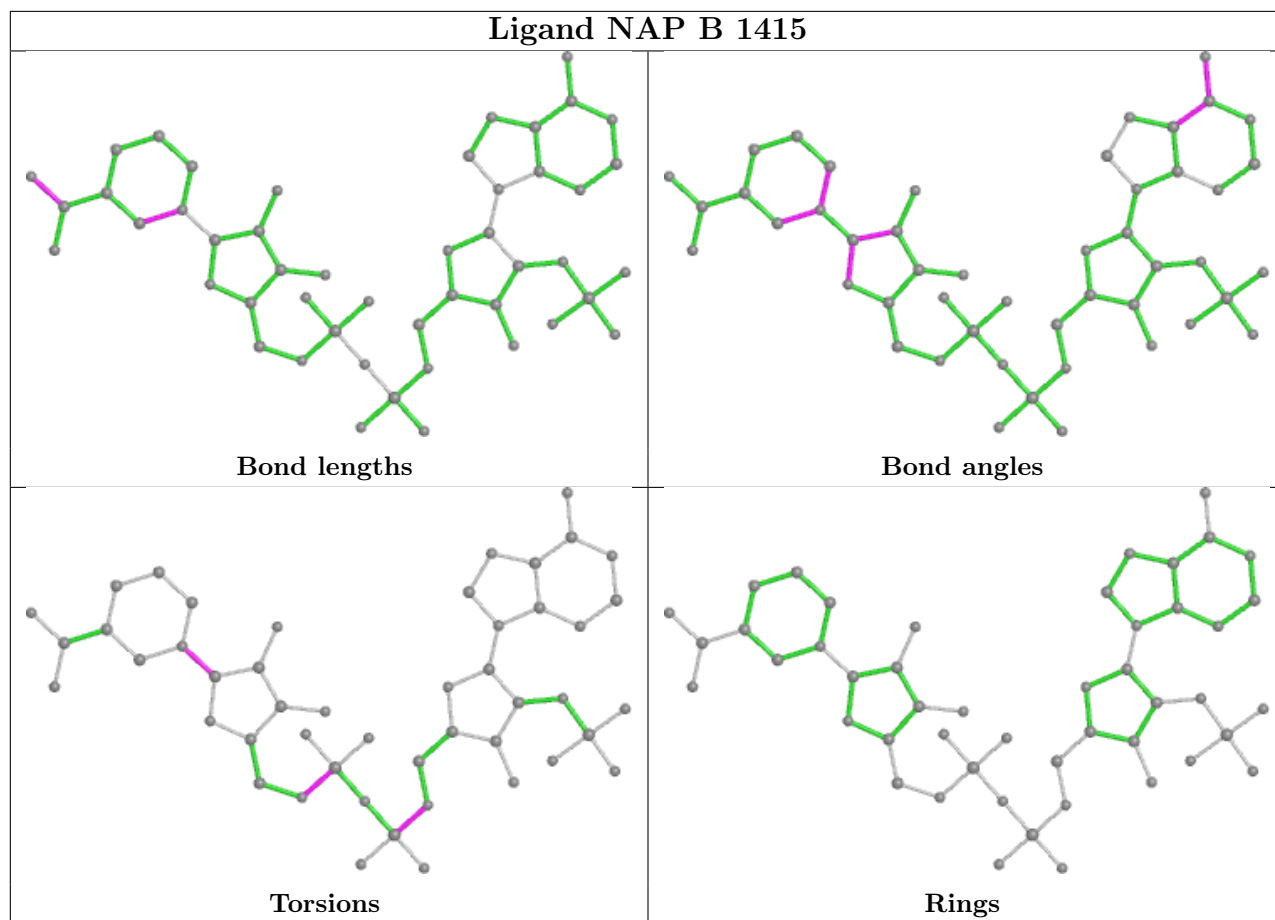
There are no ring outliers.

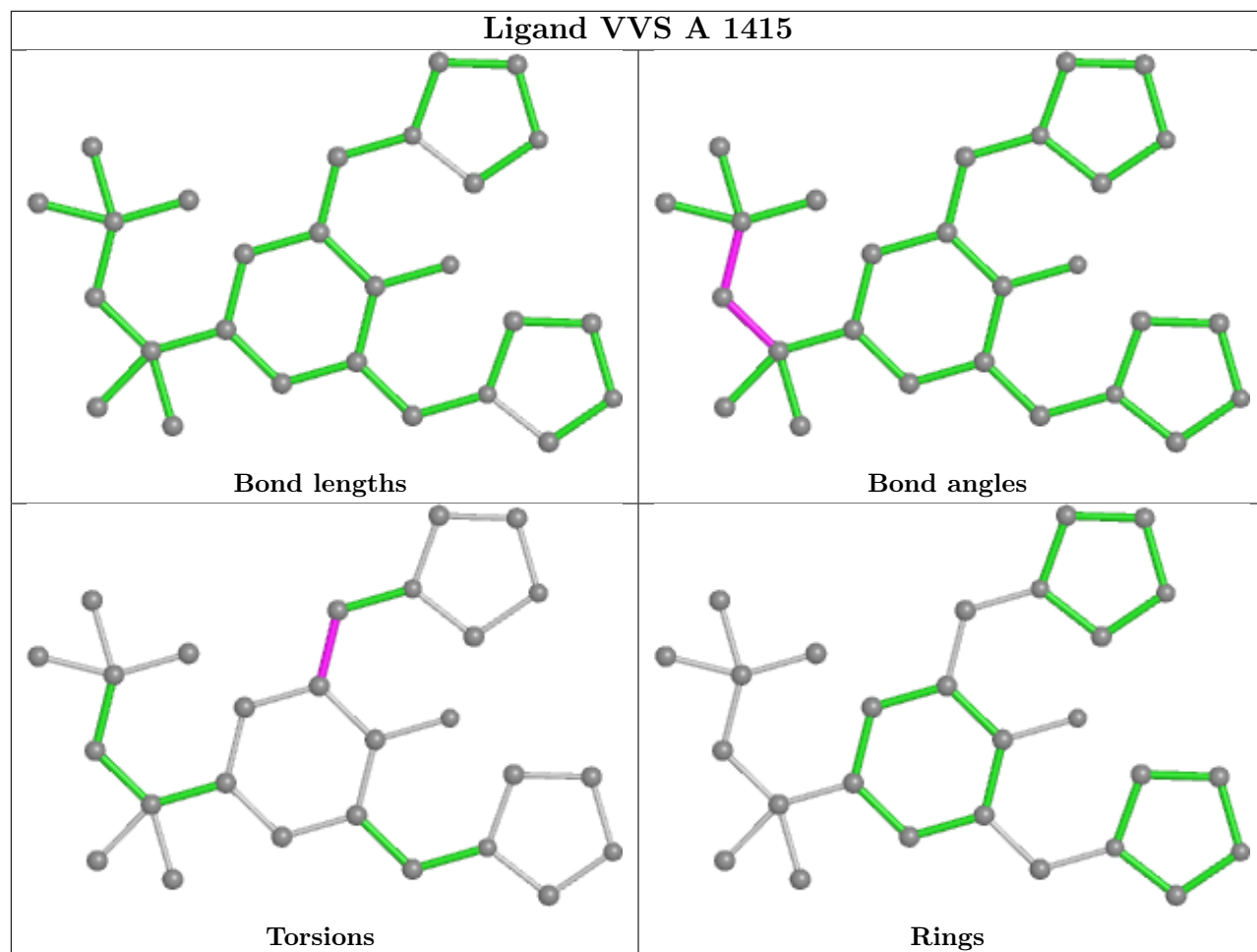
3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1416	GOL	2	0
3	A	1416	NAP	1	0
3	B	1415	NAP	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.







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	412/425 (96%)	0.34	37 (8%) 9 10	16, 27, 60, 85	0
1	B	405/425 (95%)	0.10	14 (3%) 44 45	16, 25, 55, 99	0
All	All	817/850 (96%)	0.22	51 (6%) 20 22	16, 26, 59, 99	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	286	GLY	10.1
1	B	135	TYR	8.5
1	A	284	GLY	7.7
1	A	285	TYR	7.2
1	B	277	GLN	6.9
1	B	276	VAL	5.8
1	A	272	TYR	5.6
1	A	277	GLN	5.2
1	B	414	LEU	5.1
1	A	137	ASP	5.0
1	B	136	GLY	5.0
1	A	280	SER	4.9
1	B	139	TYR	4.5
1	B	278	SER	4.2
1	A	243	LYS	4.1
1	B	272	TYR	4.0
1	A	136	GLY	3.8
1	A	251	ILE	3.8
1	B	137	ASP	3.8
1	A	252	ASP	3.8
1	A	275	ASP	3.7
1	A	283	GLN	3.7
1	A	278	SER	3.6
1	A	273	ASP	3.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	120	LEU	3.4
1	A	274	GLY	3.4
1	A	135	TYR	3.4
1	A	286	GLY	3.3
1	B	259	MET	3.2
1	A	239	PHE	3.1
1	A	237	SER	3.0
1	A	244	ILE	3.0
1	A	134	ALA	2.9
1	A	125	VAL	2.9
1	A	260	LYS	2.9
1	A	241	ALA	2.9
1	A	245	TRP	2.8
1	A	201	LEU	2.8
1	A	267	TRP	2.7
1	A	223	PHE	2.7
1	A	276	VAL	2.7
1	B	140	ARG	2.6
1	A	139	TYR	2.3
1	A	256	ALA	2.3
1	B	133	HIS	2.2
1	A	281	VAL	2.2
1	A	236	LYS	2.2
1	B	125	VAL	2.2
1	A	138	GLN	2.1
1	A	208	TYR	2.1
1	A	140	ARG	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.

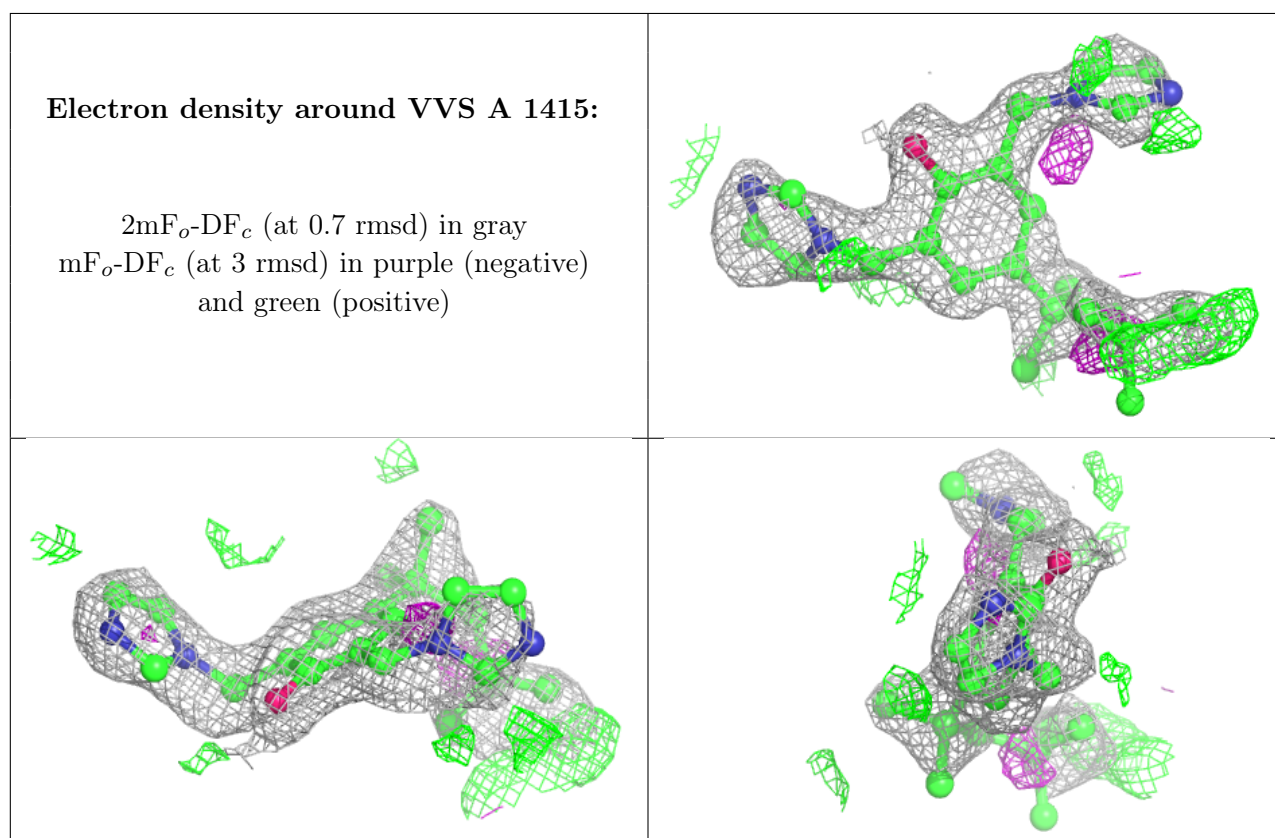
6.4 Ligands [i](#)

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, 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.

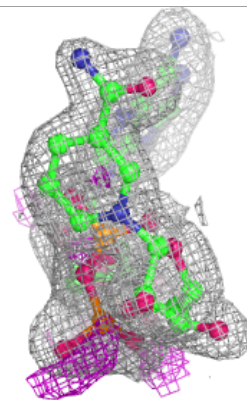
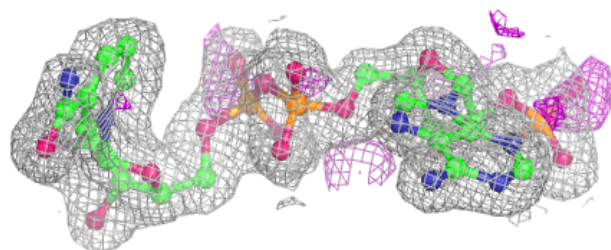
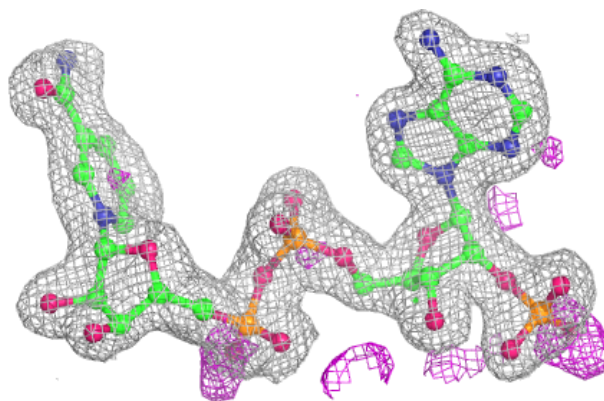
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	B	1416	6/6	0.63	0.18	57,58,59,59	0
2	VVS	A	1415	27/27	0.81	0.27	46,56,68,69	0
3	NAP	B	1415	48/48	0.96	0.10	18,28,35,40	0
3	NAP	A	1416	48/48	0.97	0.10	16,23,32,37	0

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.

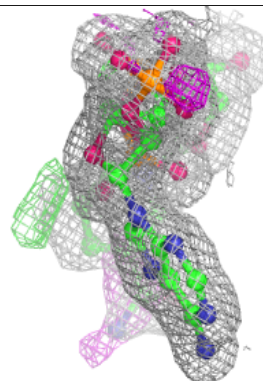
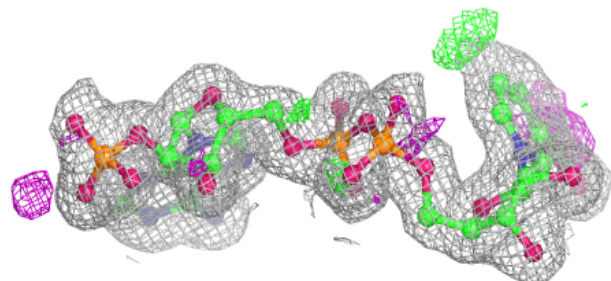
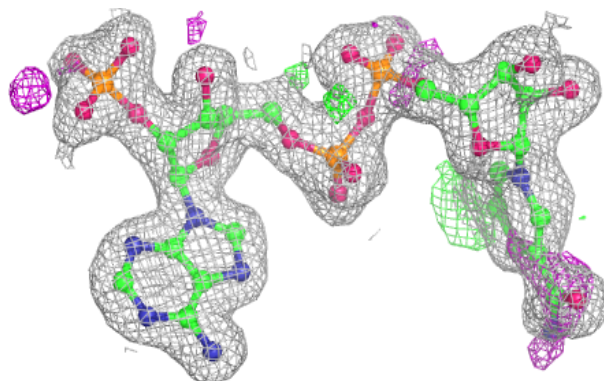


Electron density around NAP B 1415:

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

**Electron density around NAP A 1416:**

$2mF_o-DF_c$ (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.