



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 01:13 am BST

PDB ID : 4UMY
Title : IDH1 R132H in complex with cpd 1
Authors : McLean, L.; Zhang, Y.; Mathieu, M.
Deposited on : 2014-05-22
Resolution : 2.07 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

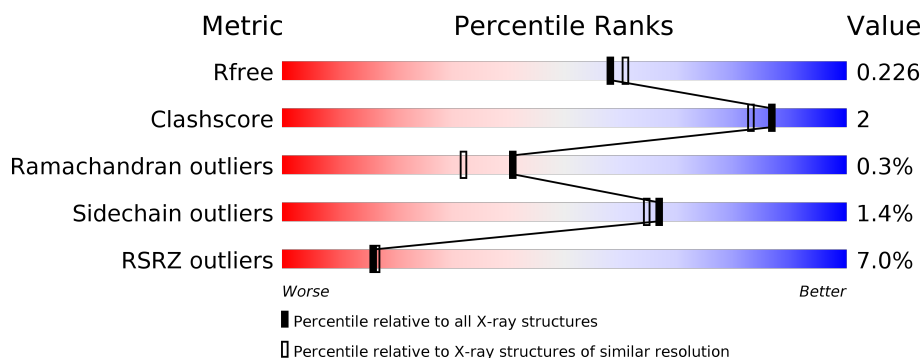
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.07 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	2684 (2.08-2.04)
Clashscore	141614	2801 (2.08-2.04)
Ramachandran outliers	138981	2768 (2.08-2.04)
Sidechain outliers	138945	2768 (2.08-2.04)
RSRZ outliers	127900	2646 (2.08-2.04)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	<div> <div>4%</div> <div> <div></div> <div>88%</div> <div>•</div> <div>8%</div> </div> </div>
1	B	425	<div> <div>8%</div> <div> <div></div> <div>83%</div> <div>7%</div> <div>10%</div> </div> </div>

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	SO4	B	1415	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6538 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
1	A	391	Total	C	N	O	S	0	0	0
			3102	1976	529	579	18			
1	B	382	Total	C	N	O	S	0	0	0
			3028	1929	516	565	18			

There are 24 discrepancies between the modelled and reference sequences:

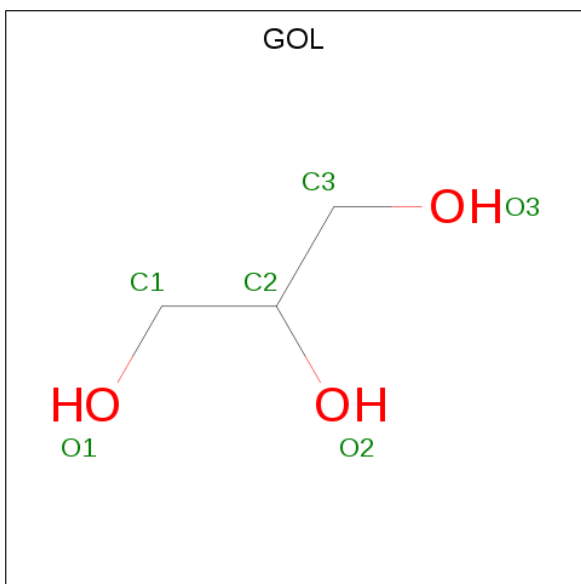
Chain	Residue	Modelled	Actual	Comment	Reference
A	415	SER	-	expression tag	UNP O75874
A	416	LEU	-	expression tag	UNP O75874
A	417	GLU	-	expression tag	UNP O75874
A	418	HIS	-	expression tag	UNP O75874
A	419	HIS	-	expression tag	UNP O75874
A	420	HIS	-	expression tag	UNP O75874
A	421	HIS	-	expression tag	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 SULFATE ION (three-letter code: SO4) (formula: O₄S).



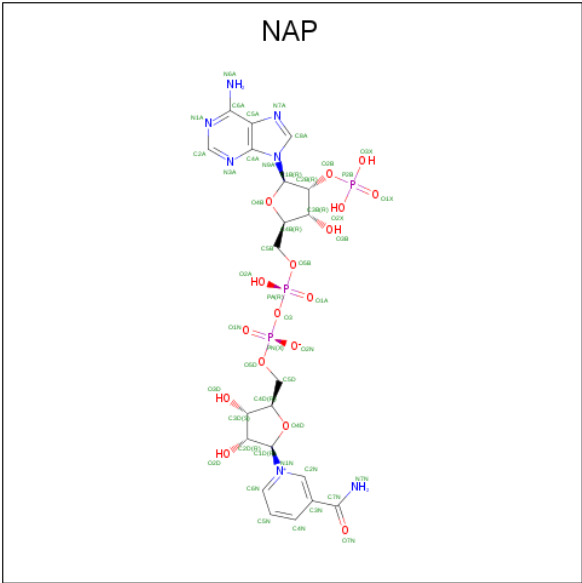
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

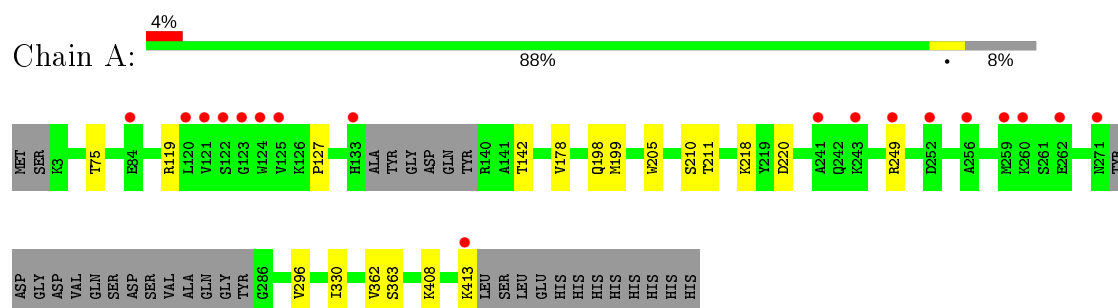
- Molecule 4 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



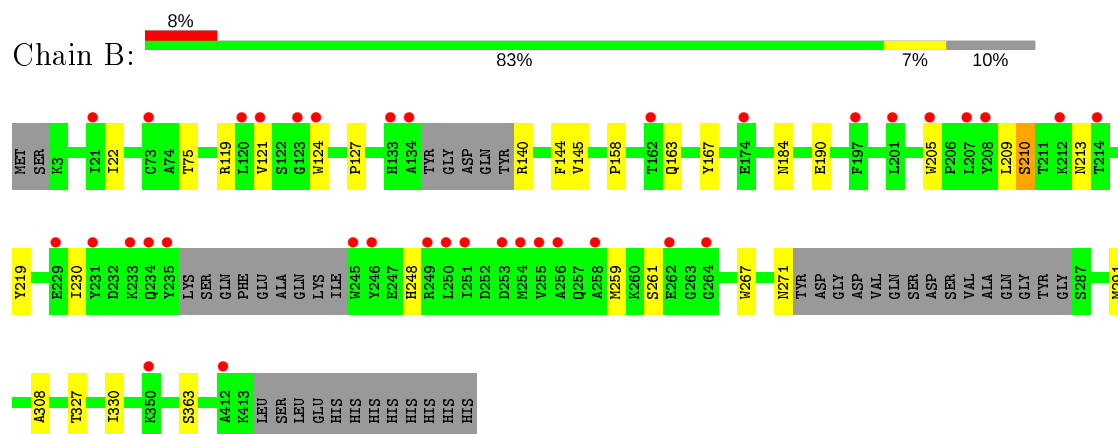
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: ISOCITRATE DEHYDROGENASE [NADP] CYTOPLASMIC



- Molecule 1: ISOCITRATE DEHYDROGENASE [NADP] CYTOPLASMIC



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	82.68 Å 82.68 Å 300.37 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.34 – 2.07 41.34 – 2.07	Depositor EDS
% Data completeness (in resolution range)	99.2 (41.34-2.07) 99.5 (41.34-2.07)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.45 (at 2.06 Å)	Xtriage
Refinement program	BUSTER 2.11.5	Depositor
R, R_{free}	0.200 , 0.224 0.198 , 0.226	Depositor DCC
R_{free} test set	3248 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	43.2	Xtriage
Anisotropy	0.243	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 56.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6538	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% 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: GOL, NAP, SO4

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	0.51	0/3165	0.62	0/4262
1	B	0.51	0/3089	0.64	0/4161
All	All	0.51	0/6254	0.63	0/8423

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	3102	0	3097	15	0
1	B	3028	0	3020	15	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
3	A	6	0	8	1	0
4	A	48	0	25	1	0
4	B	48	0	25	1	0
5	A	184	0	0	1	0
5	B	102	0	0	0	0
All	All	6538	0	6175	26	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 2.

All (26) 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:178:VAL:HG21	1:B:219:TYR:HA	1.33	1.10
1:A:218:LYS:HE2	1:B:145:VAL:HG23	1.81	0.62
1:A:199:MET:CE	1:A:296:VAL:HG21	2.29	0.62
1:A:218:LYS:HD3	1:B:144:PHE:HA	1.80	0.62
1:A:330:ILE:HD12	1:A:363:SER:HB3	1.86	0.56
1:B:291:MET:HB3	1:B:308:ALA:HB3	1.88	0.54
1:A:75:THR:O	4:A:1417:NAP:H2N	2.09	0.53
1:A:362:VAL:HG23	1:A:408:LYS:HD2	1.92	0.51
1:A:362:VAL:HG23	1:A:408:LYS:CE	2.42	0.50
1:B:330:ILE:HD12	1:B:363:SER:HB3	1.94	0.49
1:B:190:GLU:HG2	1:B:230:ILE:HD11	1.95	0.48
3:A:1416:GOL:H32	5:A:2178:HOH:O	2.12	0.48
1:B:158:PRO:HD2	1:B:163:GLN:O	2.14	0.48
1:B:75:THR:O	4:B:1416:NAP:H2N	2.14	0.47
1:A:142:THR:HG21	1:B:167:TYR:HB3	1.96	0.46
1:A:210:SER:HA	1:A:249:ARG:O	2.16	0.46
1:B:22:ILE:HD11	1:B:327:THR:HB	1.97	0.45
1:B:209:LEU:HD23	1:B:248:HIS:HD2	1.82	0.44
1:A:178:VAL:CG2	1:B:219:TYR:HA	2.25	0.44
1:A:211:THR:HB	1:A:220:ASP:HB3	2.01	0.41
1:B:121:VAL:HB	1:B:124:TRP:CE2	2.55	0.41
1:B:210:SER:HB3	1:B:267:TRP:HE1	1.85	0.41
1:A:127:PRO:HD2	1:A:205:TRP:CH2	2.55	0.41
1:A:199:MET:HE1	1:A:296:VAL:HG21	2.00	0.41
1:B:127:PRO:HD2	1:B:205:TRP:CH2	2.55	0.41
1:A:199:MET:HE2	1:A:296:VAL:HG21	1.99	0.41

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	385/425 (91%)	376 (98%)	9 (2%)	0	100	100
1	B	374/425 (88%)	361 (96%)	11 (3%)	2 (0%)	29	19
All	All	759/850 (89%)	737 (97%)	20 (3%)	2 (0%)	41	32

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	213	ASN
1	B	261	SER

5.3.2 Protein sidechains ⓘ

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	332/361 (92%)	329 (99%)	3 (1%)	78	78
1	B	324/361 (90%)	318 (98%)	6 (2%)	57	53
All	All	656/722 (91%)	647 (99%)	9 (1%)	67	64

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	ARG
1	A	198	GLN
1	A	413	LYS
1	B	119	ARG
1	B	140	ARG
1	B	184	ASN
1	B	210	SER
1	B	259	MET
1	B	271	ASN

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

Mol	Chain	Res	Type
1	A	411	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

7 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	NAP	B	1416	-	45,52,52	0.99	2 (4%)	56,80,80	0.92	4 (7%)
2	SO4	B	1414	-	4,4,4	0.26	0	6,6,6	0.16	0
2	SO4	A	1415	-	4,4,4	0.35	0	6,6,6	0.29	0
4	NAP	A	1417	-	45,52,52	1.18	3 (6%)	56,80,80	0.94	4 (7%)
2	SO4	A	1414	-	4,4,4	0.15	0	6,6,6	0.11	0
3	GOL	A	1416	-	5,5,5	0.24	0	5,5,5	0.73	0
2	SO4	B	1415	-	4,4,4	0.31	0	6,6,6	0.18	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
4	NAP	B	1416	-	-	7/31/67/67	0/5/5/5
4	NAP	A	1417	-	-	9/31/67/67	0/5/5/5
3	GOL	A	1416	-	-	2/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1417	NAP	C2N-N1N	5.50	1.41	1.35
4	B	1416	NAP	C2N-N1N	3.89	1.39	1.35
4	B	1416	NAP	O7N-C7N	3.26	1.30	1.24
4	A	1417	NAP	O7N-C7N	3.09	1.30	1.24
4	A	1417	NAP	O4D-C1D	-2.17	1.38	1.41

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1417	NAP	O4D-C1D-C2D	-2.59	103.14	106.93
4	B	1416	NAP	O4D-C1D-C2D	-2.37	103.46	106.93
4	B	1416	NAP	C2N-N1N-C1D	-2.36	113.87	119.14
4	B	1416	NAP	C5A-C6A-N6A	2.24	123.75	120.35
4	A	1417	NAP	O2B-C2B-C1B	-2.23	102.08	110.10
4	B	1416	NAP	O3B-C3B-C2B	2.22	117.46	111.17
4	A	1417	NAP	C5A-C6A-N6A	2.17	123.65	120.35
4	A	1417	NAP	C2N-N1N-C1D	-2.14	114.37	119.14

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1416	NAP	C5B-O5B-PA-O1A
4	B	1416	NAP	C5D-O5D-PN-O2N
4	B	1416	NAP	O4D-C1D-N1N-C2N
4	B	1416	NAP	O4D-C1D-N1N-C6N
4	B	1416	NAP	C2D-C1D-N1N-C2N
4	A	1417	NAP	C5D-O5D-PN-O2N
4	A	1417	NAP	O4D-C1D-N1N-C2N
4	A	1417	NAP	O4D-C1D-N1N-C6N
4	A	1417	NAP	C2D-C1D-N1N-C2N
3	A	1416	GOL	C1-C2-C3-O3
4	A	1417	NAP	PN-O3-PA-O5B
4	B	1416	NAP	C5D-O5D-PN-O3

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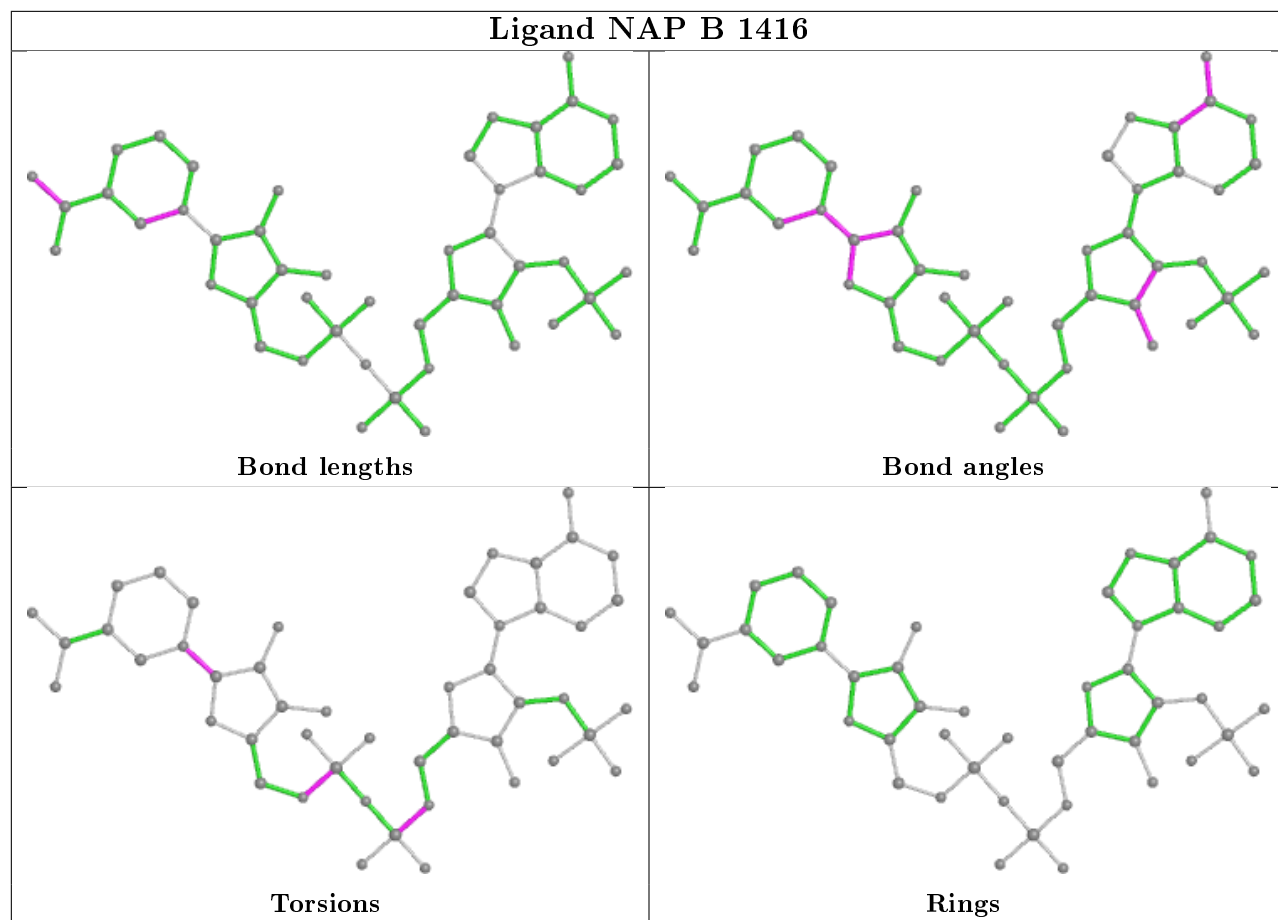
Mol	Chain	Res	Type	Atoms
4	A	1417	NAP	C5D-O5D-PN-O3
3	A	1416	GOL	O2-C2-C3-O3
4	B	1416	NAP	C5D-O5D-PN-O1N
4	A	1417	NAP	C5D-O5D-PN-O1N
4	A	1417	NAP	C2D-C1D-N1N-C6N
4	A	1417	NAP	C5B-O5B-PA-O1A

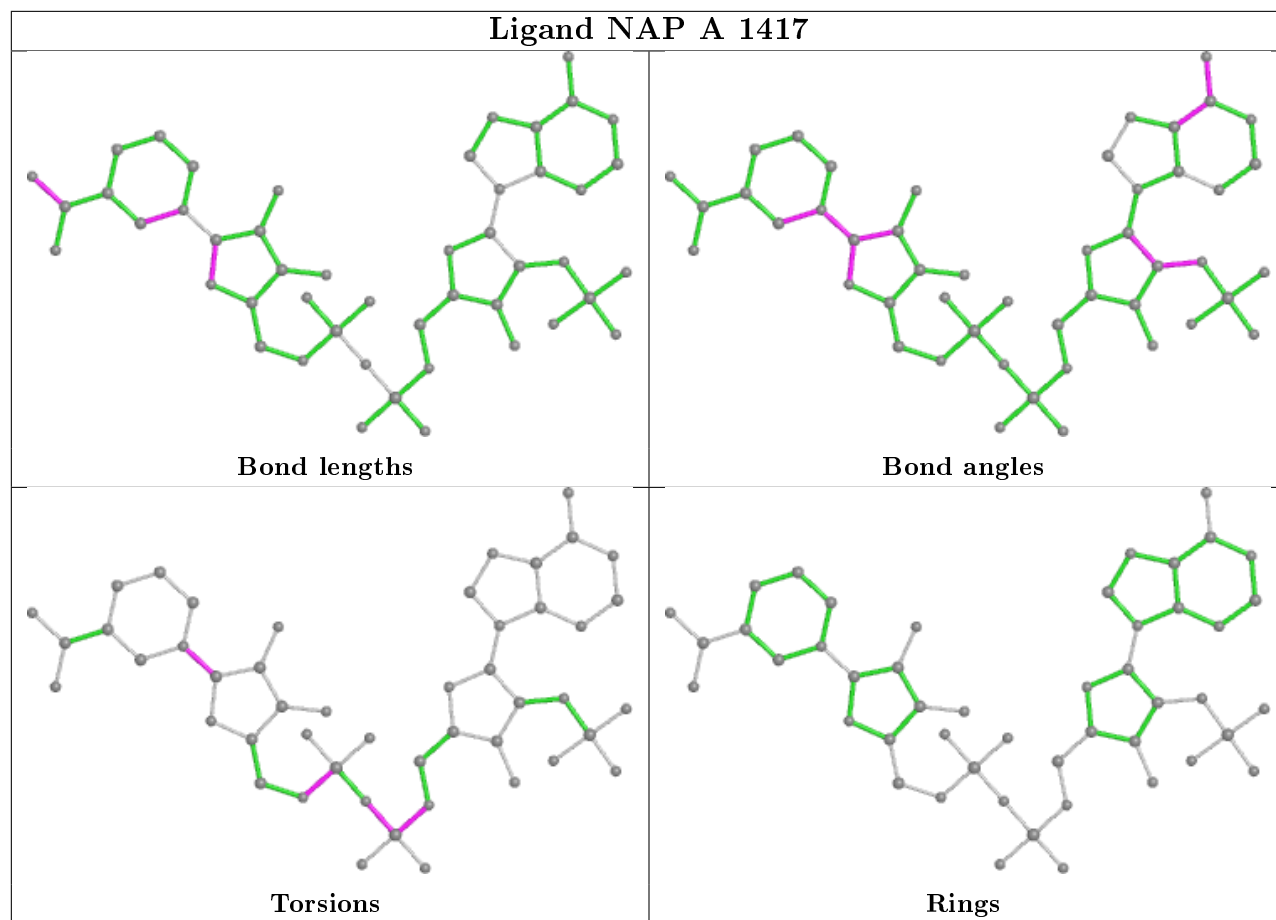
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1416	NAP	1	0
4	A	1417	NAP	1	0
3	A	1416	GOL	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	391/425 (92%)	0.34	18 (4%) 32 33	32, 44, 80, 119	0
1	B	382/425 (89%)	0.53	36 (9%) 8 9	32, 47, 93, 120	0
All	All	773/850 (90%)	0.43	54 (6%) 16 17	32, 45, 88, 120	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	201	LEU	9.2
1	A	120	LEU	6.3
1	A	123	GLY	5.1
1	B	251	ILE	4.9
1	A	125	VAL	4.5
1	B	214	THR	4.4
1	A	133	HIS	4.3
1	A	241	ALA	4.2
1	B	233	LYS	4.1
1	B	253	ASP	4.0
1	A	260	LYS	3.9
1	B	258	ALA	3.9
1	B	255	VAL	3.7
1	B	250	LEU	3.5
1	A	259	MET	3.5
1	B	245	TRP	3.5
1	B	121	VAL	3.4
1	B	412	ALA	3.4
1	B	205	TRP	3.3
1	B	256	ALA	3.1
1	B	249	ARG	3.1
1	A	124	TRP	3.0
1	B	234	GLN	3.0
1	B	124	TRP	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	208	TYR	2.8
1	B	235	TYR	2.8
1	B	246	TYR	2.8
1	A	121	VAL	2.8
1	B	21	ILE	2.7
1	A	262	GLU	2.6
1	B	264	GLY	2.6
1	B	262	GLU	2.6
1	A	243	LYS	2.5
1	A	122	SER	2.5
1	B	134	ALA	2.5
1	B	254	MET	2.5
1	B	133	HIS	2.4
1	B	123	GLY	2.4
1	B	73	CYS	2.3
1	A	271	ASN	2.3
1	B	229	GLU	2.2
1	A	249	ARG	2.2
1	A	256	ALA	2.2
1	A	413	LYS	2.2
1	B	197	PHE	2.2
1	A	84	GLU	2.2
1	B	162	THR	2.2
1	B	120	LEU	2.1
1	B	231	TYR	2.1
1	B	350	LYS	2.1
1	B	207	LEU	2.1
1	B	174	GLU	2.0
1	A	252	ASP	2.0
1	B	212	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

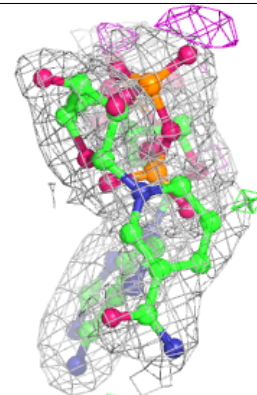
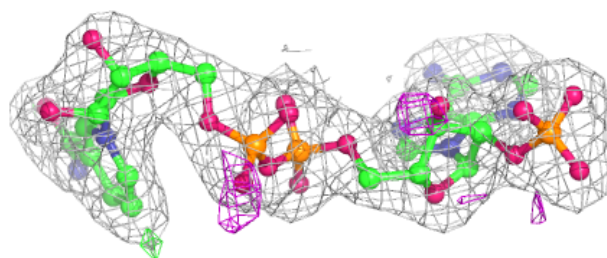
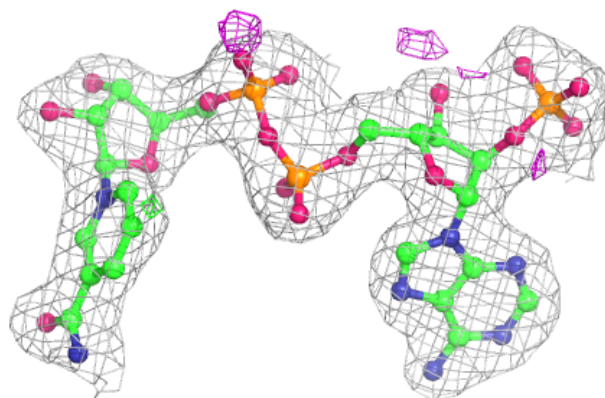
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	SO4	B	1415	5/5	0.68	0.50	114,115,116,117	0
2	SO4	A	1414	5/5	0.83	0.19	140,140,140,141	0
3	GOL	A	1416	6/6	0.86	0.22	52,56,58,59	0
2	SO4	A	1415	5/5	0.92	0.24	70,71,76,78	0
2	SO4	B	1414	5/5	0.92	0.14	102,103,103,103	0
4	NAP	A	1417	48/48	0.98	0.09	35,45,52,54	0
4	NAP	B	1416	48/48	0.98	0.10	37,43,52,54	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.

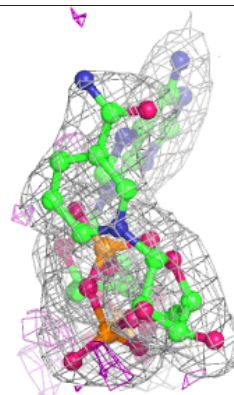
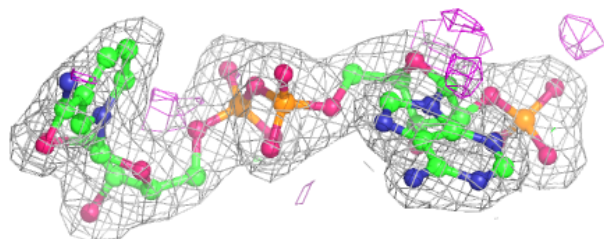
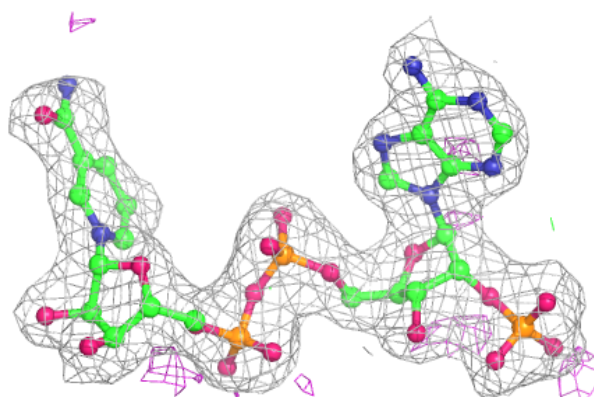
Electron density around NAP A 1417:

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 B 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.