



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:48 AM UTC

PDB ID : 4UMY / pdb_00004umy
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 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

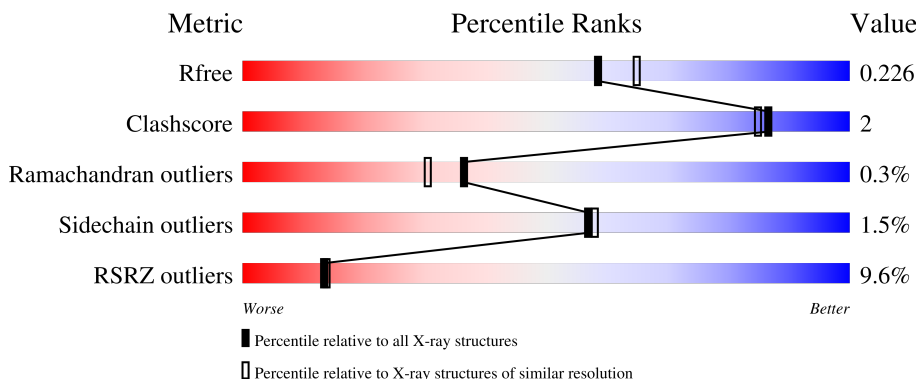
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.07 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}	180053	3774 (2.08-2.04)
Clashscore	190562	3883 (2.08-2.04)
Ramachandran outliers	187476	3860 (2.08-2.04)
Sidechain outliers	187428	3860 (2.08-2.04)
RSRZ outliers	180081	3775 (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 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	
1	B	425	

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:

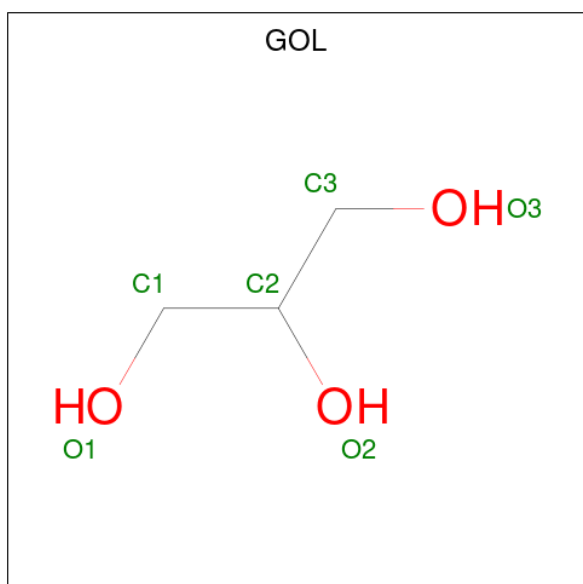
Chain	Residue	Modelled	Actual	Comment	Reference
A	415	SER	-	expression tag	UNP O75874
A	416	LEU	-	expression tag	UNP O75874
A	417	GLU	-	expression tag	UNP O75874
A	418	HIS	-	expression tag	UNP O75874
A	419	HIS	-	expression tag	UNP O75874
A	420	HIS	-	expression tag	UNP O75874
A	421	HIS	-	expression tag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



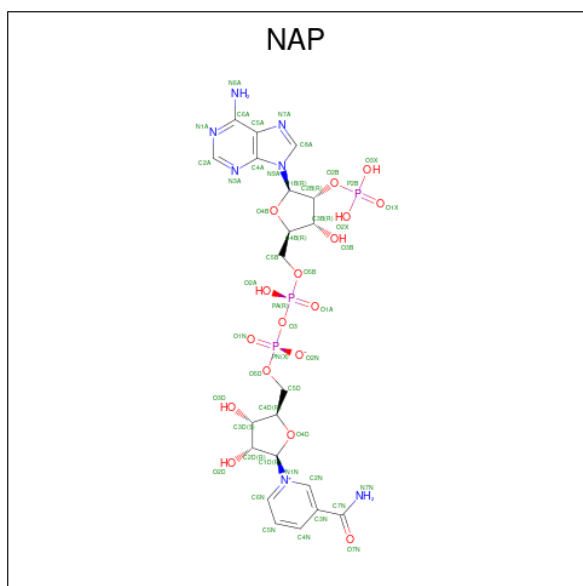
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is GLYCEROL (CCD ID: 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 (CCD ID: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



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

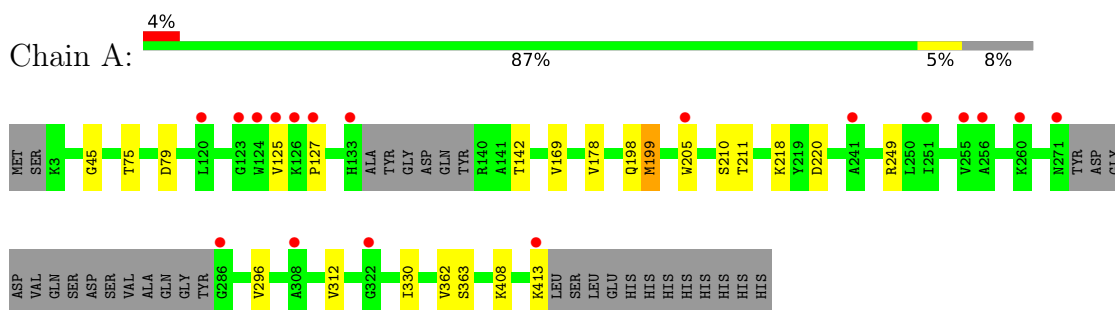
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	184	Total	O	0	0
			184	184		
5	B	102	Total	O	0	0
			102	102		

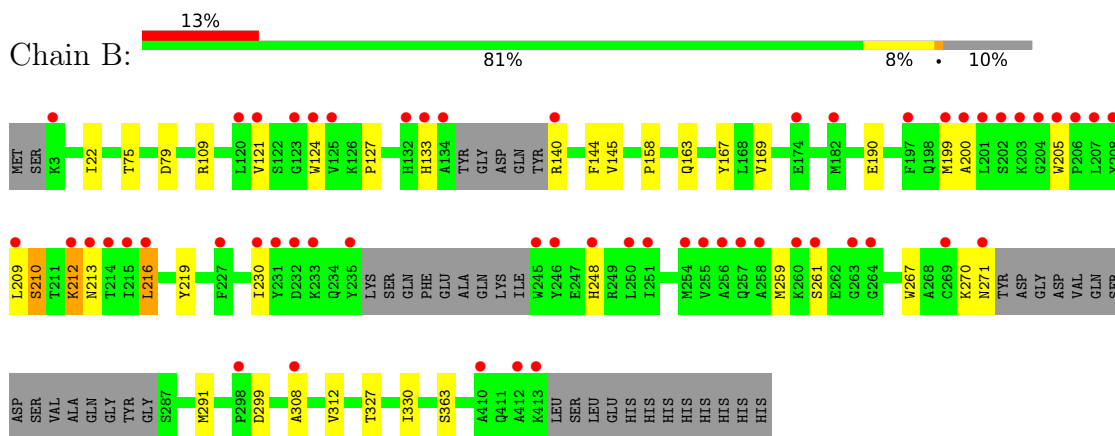
3 Residue-property plots [i](#)

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: 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	82.68Å 82.68Å 300.37Å 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 2.06Å)	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)	0.37 , 50.3	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: SO4, GOL, NAP

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	0.85	1/3165 (0.0%)	1.17	4/4262 (0.1%)
1	B	0.86	1/3089 (0.0%)	1.21	11/4161 (0.3%)
All	All	0.85	2/6254 (0.0%)	1.19	15/8423 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	199	MET	SD-CE	-8.53	1.58	1.79
1	B	133	HIS	ND1-CE1	5.39	1.38	1.32

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	200	ALA	N-CA-C	-8.45	99.72	110.19
1	B	133	HIS	CB-CG-CD2	-7.52	121.42	131.20
1	B	199	MET	N-CA-C	-6.59	103.59	111.69
1	B	212	LYS	N-CA-C	-5.83	104.56	111.03
1	A	125	VAL	N-CA-C	-5.79	106.83	111.81
1	B	169	VAL	N-CA-C	-5.62	105.37	110.82
1	B	79	ASP	CA-CB-CG	5.48	118.08	112.60
1	A	79	ASP	CA-CB-CG	5.39	117.99	112.60
1	A	169	VAL	N-CA-C	-5.27	106.55	111.45
1	A	45	GLY	N-CA-C	-5.22	105.93	111.93
1	B	270	LYS	CA-C-N	5.21	131.07	121.70
1	B	270	LYS	C-N-CA	5.21	131.07	121.70
1	B	216	LEU	CA-C-N	5.05	127.30	120.38
1	B	216	LEU	C-N-CA	5.05	127.30	120.38
1	B	299	ASP	CA-CB-CG	5.05	117.65	112.60

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	14	0
1	B	3028	0	3020	16	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:B:109:ARG:HB3	1:B:291:MET:HE1	1.76	0.66
1:A:199:MET:HE2	1:A:296:VAL:HG21	1.76	0.66
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:127:PRO:HD2	1:B:205:TRP:CH2	2.55	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%)	24	17
All	All	759/850 (89%)	737 (97%)	20 (3%)	2 (0%)	36	30

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%)	70	74
1	B	324/361 (90%)	317 (98%)	7 (2%)	45	43
All	All	656/722 (91%)	646 (98%)	10 (2%)	57	58

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	198	GLN
1	A	312	VAL
1	A	413	LYS
1	B	140	ARG
1	B	210	SER
1	B	212	LYS
1	B	216	LEU
1	B	259	MET
1	B	271	ASN
1	B	312	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	ASN
1	A	68	ASN
1	A	101	ASN
1	A	132	HIS
1	A	213	ASN
1	A	242	GLN
1	A	257	GLN
1	A	348	ASN
1	B	53	ASN
1	B	68	ASN
1	B	101	ASN
1	B	132	HIS

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Mol	Chain	Res	Type
1	B	348	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 oligosaccharides 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	$\# Z > 2$
2	SO4	B	1414	-	4,4,4	0.31	0	6,6,6	0.18	0
2	SO4	B	1415	-	4,4,4	0.41	0	6,6,6	0.22	0
4	NAP	A	1417	-	50,52,52	1.13	3 (6%)	71,80,80	0.80	3 (4%)
4	NAP	B	1416	-	50,52,52	0.91	2 (4%)	71,80,80	0.79	4 (5%)
3	GOL	A	1416	-	5,5,5	0.25	0	5,5,5	0.78	0
2	SO4	A	1415	-	4,4,4	0.44	0	6,6,6	0.28	0
2	SO4	A	1414	-	4,4,4	0.26	0	6,6,6	0.13	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
3	GOL	A	1416	-	-	2/4/4/4	-
4	NAP	A	1417	-	-	9/35/67/67	0/5/5/5
4	NAP	B	1416	-	-	7/35/67/67	0/5/5/5

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1417	NAP	C2N-N1N	6.07	1.41	1.35
4	B	1416	NAP	C2N-N1N	4.29	1.39	1.35
4	B	1416	NAP	O7N-C7N	3.37	1.30	1.24
4	A	1417	NAP	O7N-C7N	3.19	1.30	1.24
4	A	1417	NAP	O4D-C1D	-2.17	1.38	1.40

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1416	NAP	C6N-N1N-C1D	2.83	125.28	119.73
4	A	1417	NAP	C6N-N1N-C1D	2.61	124.84	119.73
4	B	1416	NAP	C2N-N1N-C1D	-2.38	113.87	119.13
4	A	1417	NAP	O2B-C2B-C1B	-2.27	102.08	110.05
4	B	1416	NAP	O3B-C3B-C2B	2.24	117.46	111.19
4	A	1417	NAP	C2N-N1N-C1D	-2.16	114.37	119.13
4	B	1416	NAP	C4D-O4D-C1D	2.05	111.80	109.92

There are no chirality outliers.

All (18) torsion outliers are listed below:

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

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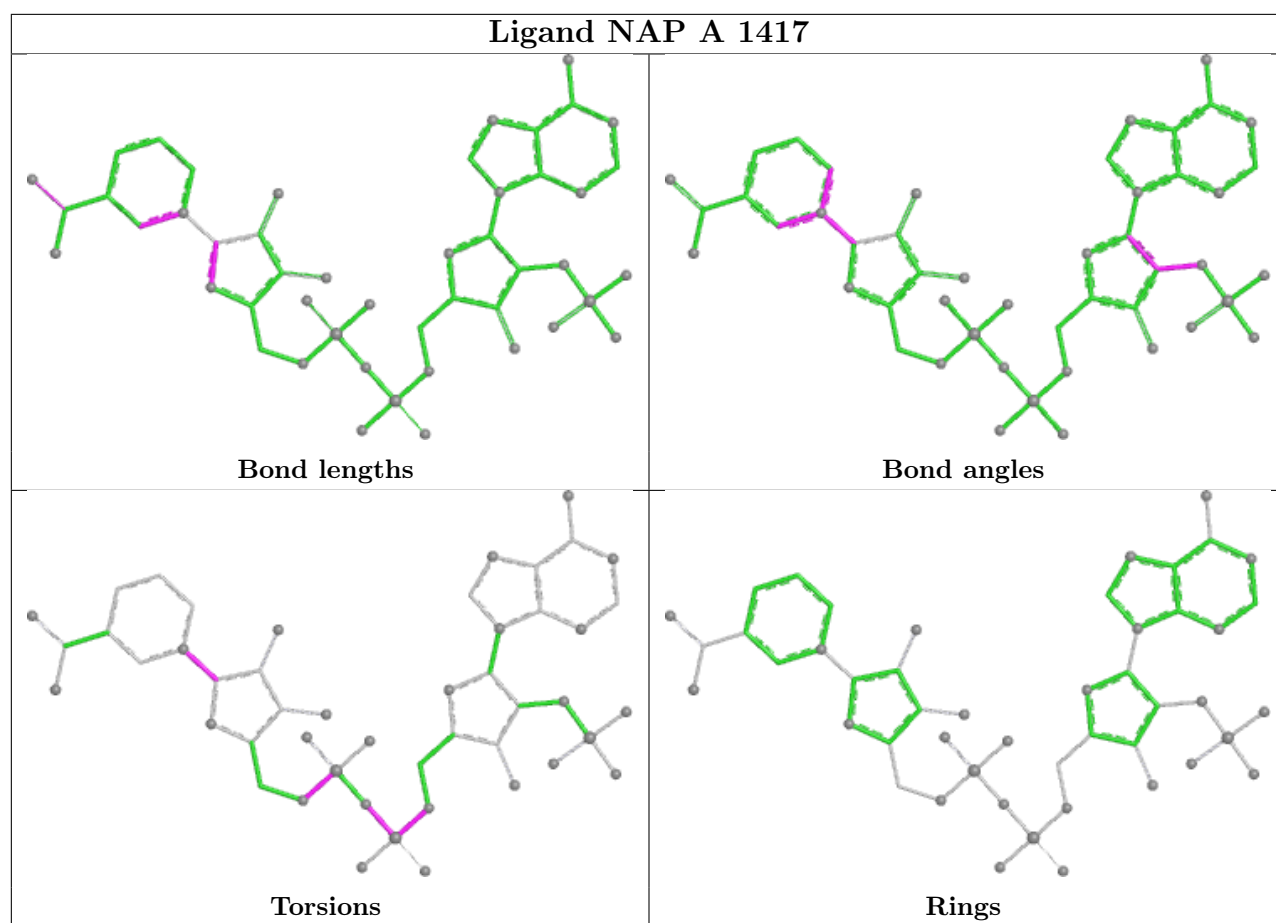
Mol	Chain	Res	Type	Atoms
3	A	1416	GOL	C1-C2-C3-O3
4	A	1417	NAP	C5B-O5B-PA-O1A
4	A	1417	NAP	C5D-O5D-PN-O1N
4	B	1416	NAP	C5D-O5D-PN-O1N
4	A	1417	NAP	C2D-C1D-N1N-C6N

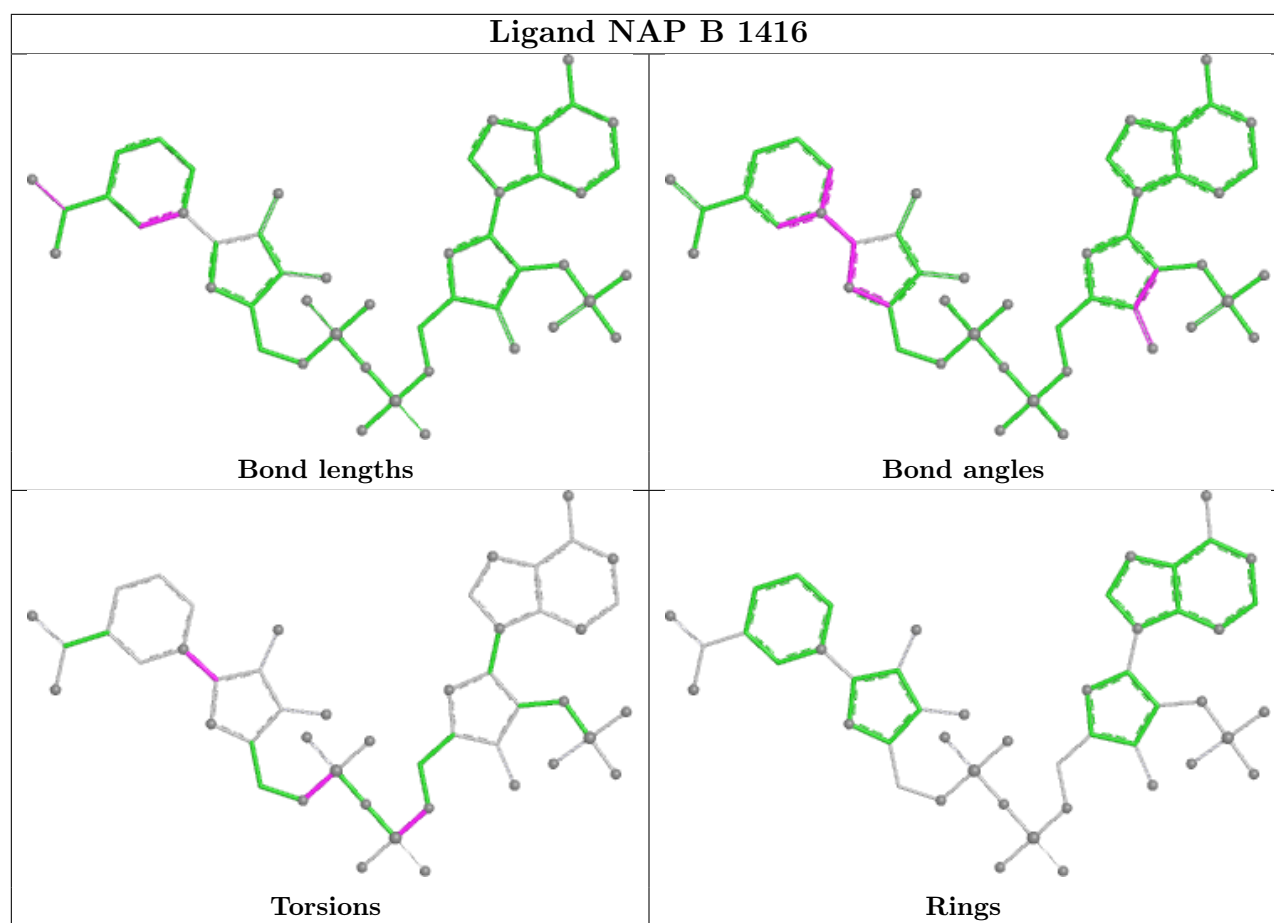
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1417	NAP	1	0
4	B	1416	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%)	0.27	18 (4%) 37 37	32, 44, 80, 119	0
1	B	382/425 (89%)	0.66	56 (14%) 6 5	32, 47, 93, 120	0
All	All	773/850 (90%)	0.46	74 (9%) 13 14	32, 45, 88, 120	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	245	TRP	7.2
1	B	201	LEU	6.6
1	B	251	ILE	6.1
1	A	120	LEU	6.1
1	A	133	HIS	4.8
1	B	235	TYR	4.6
1	B	125	VAL	4.5
1	B	255	VAL	4.4
1	A	125	VAL	4.2
1	B	134	ALA	4.1
1	B	258	ALA	4.1
1	B	205	TRP	4.1
1	B	200	ALA	4.1
1	B	215	ILE	4.0
1	B	204	GLY	3.9
1	B	250	LEU	3.8
1	B	208	TYR	3.8
1	B	133	HIS	3.7
1	B	197	PHE	3.6
1	B	132	HIS	3.5
1	B	230	ILE	3.5
1	B	121	VAL	3.4
1	B	233	LYS	3.4
1	B	207	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	231	TYR	3.2
1	B	246	TYR	3.2
1	B	206	PRO	3.2
1	B	413	LYS	3.2
1	B	264	GLY	3.1
1	B	124	TRP	3.0
1	A	241	ALA	3.0
1	B	216	LEU	2.9
1	B	214	THR	2.9
1	A	271	ASN	2.9
1	A	123	GLY	2.8
1	A	413	LYS	2.8
1	A	286	GLY	2.8
1	A	251	ILE	2.8
1	A	127	PRO	2.7
1	B	203	LYS	2.7
1	B	412	ALA	2.7
1	A	205	TRP	2.7
1	B	257	GLN	2.6
1	B	120	LEU	2.6
1	B	182	MET	2.6
1	B	232	ASP	2.5
1	B	209	LEU	2.5
1	A	322	GLY	2.5
1	B	263	GLY	2.5
1	B	308	ALA	2.4
1	B	248	HIS	2.4
1	B	271	ASN	2.4
1	A	308	ALA	2.4
1	B	140	ARG	2.4
1	B	202	SER	2.4
1	A	256	ALA	2.4
1	B	227	PHE	2.3
1	B	254	MET	2.3
1	B	269	CYS	2.3
1	A	124	TRP	2.3
1	B	213	ASN	2.3
1	B	123	GLY	2.3
1	A	126	LYS	2.2
1	A	255	VAL	2.2
1	B	3	LYS	2.2
1	B	298	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	256	ALA	2.2
1	B	261	SER	2.2
1	B	199	MET	2.2
1	B	174	GLU	2.1
1	B	410	ALA	2.1
1	A	260	LYS	2.1
1	B	212	LYS	2.1
1	B	260	LYS	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 [i](#)

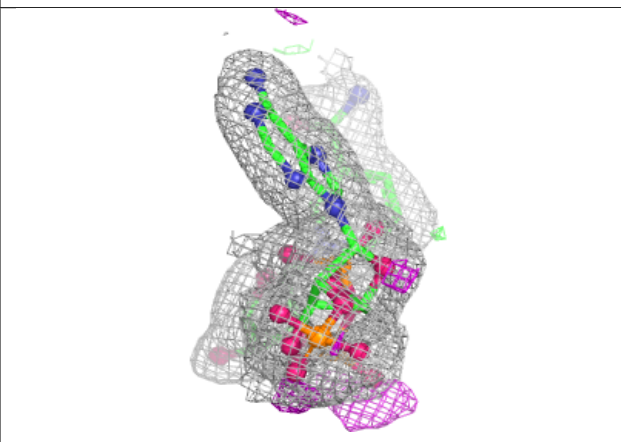
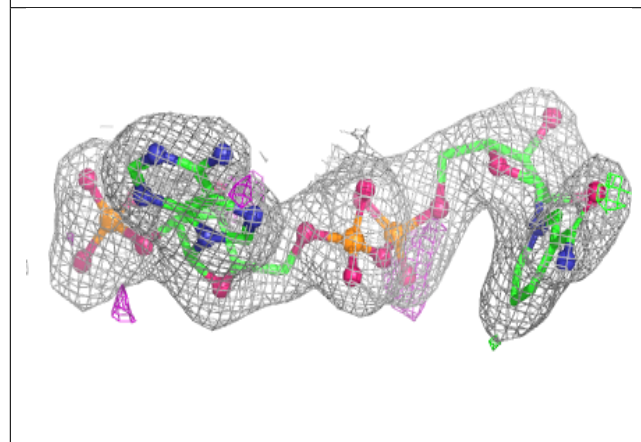
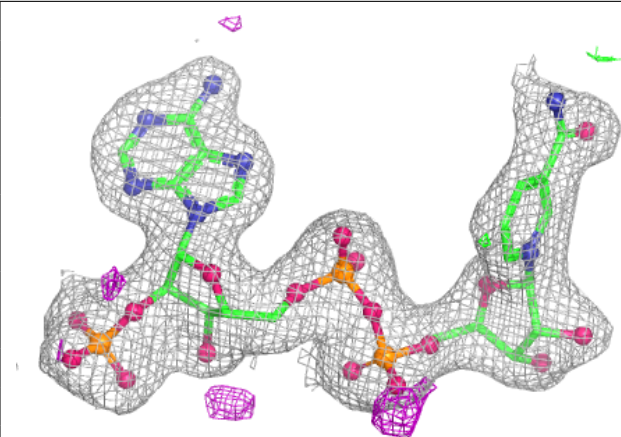
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	B	1415	5/5	0.71	0.20	114,115,116,117	0
2	SO4	A	1414	5/5	0.76	0.16	140,140,140,141	0
2	SO4	B	1414	5/5	0.78	0.13	102,103,103,103	0
3	GOL	A	1416	6/6	0.79	0.15	52,56,58,59	0
2	SO4	A	1415	5/5	0.86	0.11	70,71,76,78	0
4	NAP	A	1417	48/48	0.96	0.07	35,45,52,54	0
4	NAP	B	1416	48/48	0.96	0.07	37,43,52,54	0

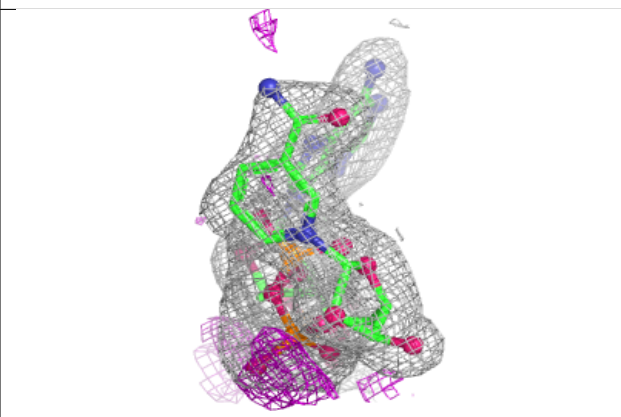
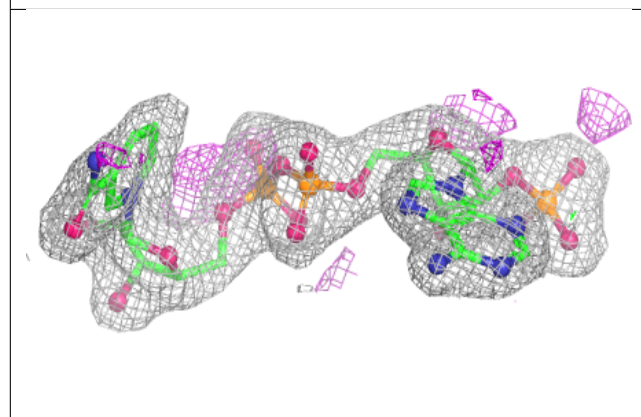
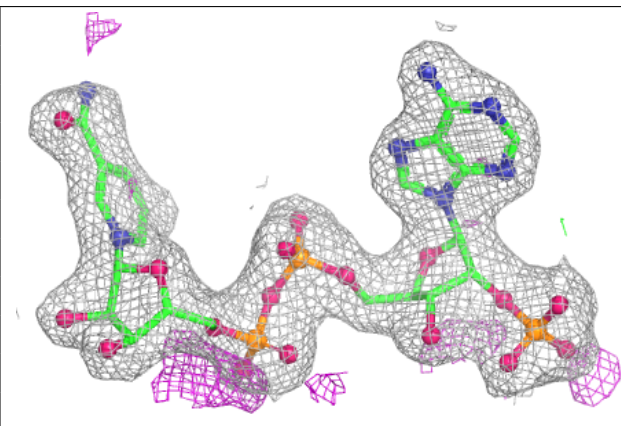
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation 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.