



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 07:13 PM BST

PDB ID : 5V7N
Title : Crystal structure of NADPH-dependent glyoxylate/hydroxypyruvate reductase SMC04462 (SmGhrB) from *Sinorhizobium meliloti* in complex with NADP and 2-Keto-D-gluconic acid
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Deposited on : 2017-03-20
Resolution : 1.75 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	:	2.13.1
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)	:	2.13.1

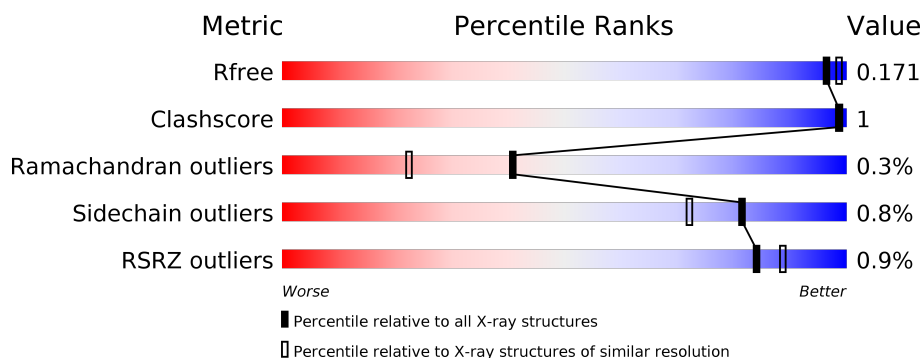
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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	323	<div> <div>%</div> <div> <div></div> <div>97%</div> <div>..</div> </div> </div>
1	B	323	<div> <div>%</div> <div> <div></div> <div>96%</div> <div>..</div> </div> </div>
1	C	323	<div> <div>%</div> <div> <div></div> <div>95%</div> <div>..</div> </div> </div>
1	D	323	<div> <div>%</div> <div> <div></div> <div>97%</div> <div>..</div> </div> </div>

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 11162 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.

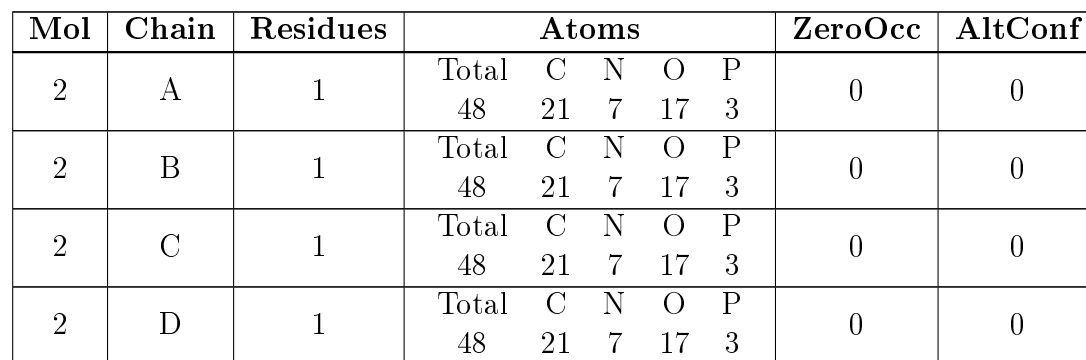
- Molecule 1 is a protein called NADPH-dependent glyoxylate/hydroxypyruvate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	S	0	4	0
			2396	1522	427	442	5			
1	B	318	Total	C	N	O	S	0	1	0
			2364	1503	418	438	5			
1	C	318	Total	C	N	O	S	0	4	0
			2377	1515	419	438	5			
1	D	318	Total	C	N	O	S	0	1	0
			2374	1511	421	437	5			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q92LZ4
B	0	SER	-	expression tag	UNP Q92LZ4
C	0	SER	-	expression tag	UNP Q92LZ4
D	0	SER	-	expression tag	UNP Q92LZ4

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



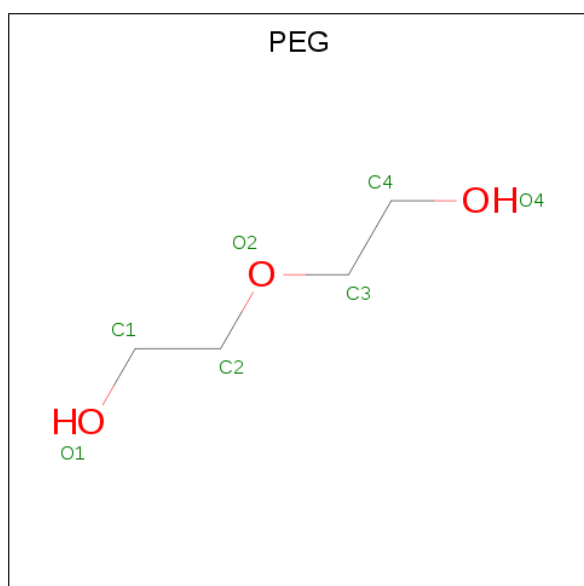
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- Chemical structure of 8YV (8-epi-yunnanin) is shown, featuring a central carbon atom (C3(S)) bonded to a hydroxyl group (O4), a hydroxyl group (O3), a carboxylic acid group (C1=O2, O1A), and a side chain (C4(R), C5(R), C6, O6). The structure is labeled with atom names (C1, C2, C3(S), C4(R), C5(R), C6, O1A, O2, O3, O4, O5, O6) and stereochemistry (R, S).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	6	7		
3	B	1	Total	C	O	0	0
			13	6	7		
3	C	1	Total	C	O	0	0
			13	6	7		
3	D	1	Total	C	O	0	0
			13	6	7		

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Na	0	0
			1	1		
4	A	1	Total	Na	0	0
			1	1		
4	D	1	Total	Na	0	0
			1	1		
4	C	1	Total	Na	0	0
			1	1		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).

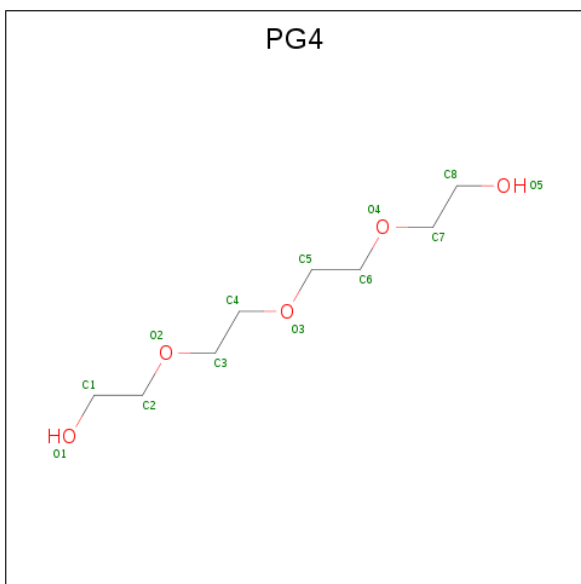


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		
5	D	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	2	Total	Cl	0	0
			2	2		
6	A	2	Total	Cl	0	0
			2	2		
6	D	2	Total	Cl	0	0
			2	2		
6	C	1	Total	Cl	0	0
			1	1		

- Molecule 7 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C₈H₁₈O₅).

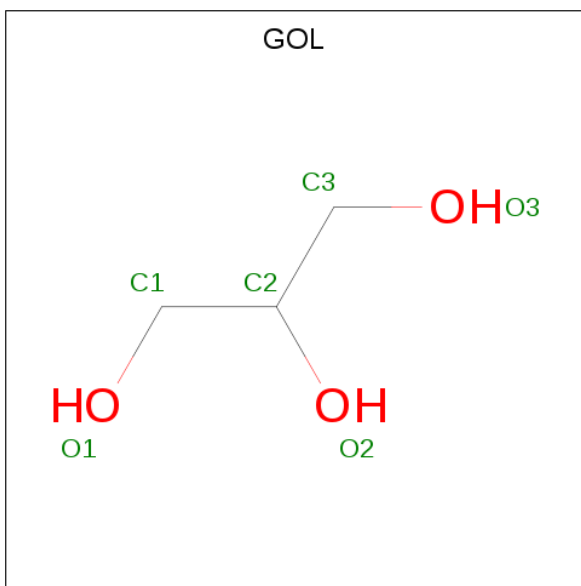


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			13	8	5		

- Molecule 8 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	2	Total	Mg	0	0
			2	2		
8	A	2	Total	Mg	0	0
			2	2		
8	D	2	Total	Mg	0	0
			2	2		
8	C	1	Total	Mg	0	0
			1	1		

- Molecule 9 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

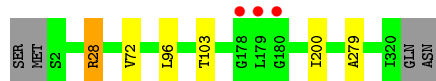
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	367	Total	O	0	2
			368	368		
10	B	360	Total	O	0	1
			361	361		
10	C	312	Total	O	0	2
			314	314		
10	D	307	Total	O	0	1
			307	307		

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: NADPH-dependent glyoxylate/hydroxypyruvate reductase



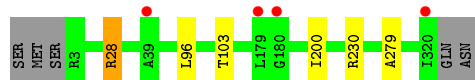
- Molecule 1: NADPH-dependent glyoxylate/hydroxypyruvate reductase



- Molecule 1: NADPH-dependent glyoxylate/hydroxypyruvate reductase



- Molecule 1: NADPH-dependent glyoxylate/hydroxypyruvate reductase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	175.90Å 175.90Å 135.34Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.11 – 1.75 45.11 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.4 (45.11-1.75) 99.3 (45.11-1.75)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.144 , 0.167 0.151 , 0.171	Depositor DCC
R_{free} test set	7842 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 34.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.237 for h,-h-k,-l	Xtriage
Reported twinning fraction	0.247 for H, K, L 0.753 for K, H, -L	Depositor
Outliers	0 of 156732 reflections	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	11162	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 48.10 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 9.0298e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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, MG, 8YV, CL, NA, NAP, PG4, PEG

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	0.61	0/2455	0.87	6/3353 (0.2%)
1	B	0.61	0/2410	0.81	5/3295 (0.2%)
1	C	0.54	0/2434	0.83	11/3327 (0.3%)
1	D	0.51	0/2420	0.81	5/3307 (0.2%)
All	All	0.57	0/9719	0.83	27/13282 (0.2%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	28[A]	ARG	NE-CZ-NH1	11.26	125.93	120.30
1	A	28[B]	ARG	NE-CZ-NH1	11.26	125.93	120.30
1	D	28	ARG	NE-CZ-NH1	9.64	125.12	120.30
1	B	28	ARG	NE-CZ-NH1	9.45	125.03	120.30
1	C	28[A]	ARG	NE-CZ-NH1	8.72	124.66	120.30
1	C	28[B]	ARG	NE-CZ-NH1	8.72	124.66	120.30
1	D	28	ARG	NE-CZ-NH2	-7.29	116.65	120.30
1	B	28	ARG	NE-CZ-NH2	-7.08	116.76	120.30
1	B	5	ARG	NE-CZ-NH1	6.69	123.65	120.30
1	C	28[A]	ARG	NE-CZ-NH2	-6.47	117.06	120.30
1	C	28[B]	ARG	NE-CZ-NH2	-6.47	117.06	120.30
1	A	28[A]	ARG	NE-CZ-NH2	-6.42	117.09	120.30
1	A	28[B]	ARG	NE-CZ-NH2	-6.42	117.09	120.30
1	C	5	ARG	NE-CZ-NH1	6.32	123.46	120.30
1	C	113	ARG	NE-CZ-NH2	-6.18	117.21	120.30
1	B	113	ARG	NE-CZ-NH2	-5.79	117.40	120.30
1	C	141	ARG	NE-CZ-NH1	5.57	123.09	120.30
1	D	28	ARG	CD-NE-CZ	5.52	131.33	123.60
1	D	230	ARG	NE-CZ-NH1	5.51	123.06	120.30
1	C	28[A]	ARG	CD-NE-CZ	5.48	131.27	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	28[B]	ARG	CD-NE-CZ	5.48	131.27	123.60
1	B	28	ARG	CD-NE-CZ	5.44	131.22	123.60
1	C	316	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	C	143	ARG	NE-CZ-NH2	-5.07	117.76	120.30
1	A	28[A]	ARG	CD-NE-CZ	5.06	130.69	123.60
1	A	28[B]	ARG	CD-NE-CZ	5.06	130.69	123.60
1	D	28	ARG	CG-CD-NE	5.05	122.41	111.80

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	2396	0	2473	2	0
1	B	2364	0	2419	1	0
1	C	2377	0	2441	2	0
1	D	2374	0	2446	1	0
2	A	48	0	25	1	0
2	B	48	0	25	0	0
2	C	48	0	25	0	0
2	D	48	0	25	0	0
3	A	13	0	0	1	0
3	B	13	0	0	1	0
3	C	13	0	0	1	0
3	D	13	0	0	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	7	0	10	0	0
5	D	7	0	10	0	0
6	A	2	0	0	0	0
6	B	2	0	0	0	0
6	C	1	0	0	0	0
6	D	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	13	0	18	0	0
8	A	2	0	0	0	0
8	B	2	0	0	0	0
8	C	1	0	0	0	0
8	D	2	0	0	0	0
9	B	6	0	8	0	0
9	D	6	0	8	0	0
10	A	368	0	0	0	0
10	B	361	0	0	0	0
10	C	314	0	0	0	0
10	D	307	0	0	0	0
All	All	11162	0	9933	10	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 1.

All (10) 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)
3:B:402:8YV:C2	3:B:402:8YV:O5	2.40	0.70
3:A:402:8YV:C2	3:A:402:8YV:O5	2.40	0.68
3:D:402:8YV:O5	3:D:402:8YV:C2	2.45	0.64
3:C:402:8YV:O5	3:C:402:8YV:C2	2.50	0.59
1:D:103:THR:HG21	1:D:279:ALA:HA	1.95	0.48
1:A:103:THR:HG21	1:A:279:ALA:HA	1.96	0.48
1:B:103:THR:HG21	1:B:279:ALA:HA	1.96	0.48
1:C:103:THR:HG21	1:C:279:ALA:HA	1.97	0.46
1:A:72:VAL:CG1	2:A:401:NAP:H6N	2.48	0.43
1:C:144[B]:THR:HG23	1:C:167[B]:SER:HB3	2.02	0.41

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	321/323 (99%)	312 (97%)	8 (2%)	1 (0%)	41	22
1	B	317/323 (98%)	309 (98%)	7 (2%)	1 (0%)	41	22
1	C	320/323 (99%)	312 (98%)	7 (2%)	1 (0%)	41	22
1	D	317/323 (98%)	308 (97%)	8 (2%)	1 (0%)	41	22
All	All	1275/1292 (99%)	1241 (97%)	30 (2%)	4 (0%)	41	22

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	200	ILE
1	A	200	ILE
1	B	200	ILE
1	D	200	ILE

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	256/258 (99%)	253 (99%)	3 (1%)	71	56
1	B	249/258 (96%)	247 (99%)	2 (1%)	81	72
1	C	250/258 (97%)	247 (99%)	3 (1%)	71	56
1	D	251/258 (97%)	249 (99%)	2 (1%)	81	72
All	All	1006/1032 (98%)	996 (99%)	10 (1%)	81	63

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

Mol	Chain	Res	Type
1	A	28[A]	ARG
1	A	28[B]	ARG
1	A	96	LEU
1	B	28	ARG

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Mol	Chain	Res	Type
1	B	96	LEU
1	C	28[A]	ARG
1	C	28[B]	ARG
1	C	96	LEU
1	D	28	ARG
1	D	96	LEU

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

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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 31 ligands modelled in this entry, 18 are monoatomic - leaving 13 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	$\# Z > 2$
7	PG4	A	407	-	12,12,12	0.49	0	11,11,11	0.38	0
9	GOL	D	407	-	5,5,5	0.49	0	5,5,5	0.61	0
5	PEG	A	404	-	6,6,6	0.45	0	5,5,5	0.17	0
2	NAP	A	401	8	45,52,52	0.85	2 (4%)	56,80,80	0.72	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAP	C	401	8	45,52,52	0.78	1 (2%)	56,80,80	0.83	0
9	GOL	B	406	-	5,5,5	0.40	0	5,5,5	0.41	0
3	8YV	B	402	-	9,12,12	0.27	0	10,16,16	1.29	1 (10%)
3	8YV	A	402	-	9,12,12	0.32	0	10,16,16	1.08	1 (10%)
3	8YV	C	402	-	9,12,12	0.19	0	10,16,16	0.74	0
2	NAP	B	401	8	45,52,52	0.87	2 (4%)	56,80,80	0.84	2 (3%)
3	8YV	D	402	-	9,12,12	0.16	0	10,16,16	0.90	0
2	NAP	D	401	8	45,52,52	0.81	2 (4%)	56,80,80	0.80	1 (1%)
5	PEG	D	404	-	6,6,6	0.46	0	5,5,5	0.15	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	PG4	A	407	-	-	1/10/10/10	-
9	GOL	D	407	-	-	4/4/4/4	-
5	PEG	A	404	-	-	3/4/4/4	-
2	NAP	A	401	8	-	2/31/67/67	0/5/5/5
2	NAP	C	401	8	-	3/31/67/67	0/5/5/5
9	GOL	B	406	-	-	0/4/4/4	-
3	8YV	B	402	-	-	0/14/18/18	-
3	8YV	A	402	-	-	1/14/18/18	-
3	8YV	C	402	-	-	4/14/18/18	-
2	NAP	B	401	8	-	3/31/67/67	0/5/5/5
3	8YV	D	402	-	-	1/14/18/18	-
2	NAP	D	401	8	-	2/31/67/67	0/5/5/5
5	PEG	D	404	-	-	1/4/4/4	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	NAP	C2A-N3A	2.21	1.35	1.32
2	C	401	NAP	PN-O2N	-2.11	1.45	1.55
2	D	401	NAP	O4B-C1B	2.09	1.44	1.41
2	D	401	NAP	PN-O2N	-2.09	1.45	1.55
2	B	401	NAP	C2A-N3A	2.07	1.35	1.32
2	A	401	NAP	PN-O2N	-2.05	1.45	1.55
2	B	401	NAP	PN-O2N	-2.01	1.45	1.55

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	8YV	O2-C2-C3	2.92	121.10	118.10
2	D	401	NAP	O4B-C1B-C2B	-2.43	102.38	106.59
3	A	402	8YV	O2-C2-C3	2.40	120.56	118.10
2	B	401	NAP	O4D-C4D-C3D	-2.17	100.82	105.11
2	B	401	NAP	O2A-PA-O1A	2.04	122.33	112.24

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	NAP	O4D-C1D-N1N-C6N
2	D	401	NAP	O4D-C1D-N1N-C6N
2	C	401	NAP	C2B-O2B-P2B-O2X
2	B	401	NAP	C2B-O2B-P2B-O1X
2	B	401	NAP	C2B-O2B-P2B-O2X
5	D	404	PEG	O2-C3-C4-O4
9	D	407	GOL	O1-C1-C2-C3
9	D	407	GOL	C1-C2-C3-O3
9	D	407	GOL	O2-C2-C3-O3
3	C	402	8YV	C3-C4-C5-O5
9	D	407	GOL	O1-C1-C2-O2
2	C	401	NAP	C2B-O2B-P2B-O1X
5	A	404	PEG	O1-C1-C2-O2
5	A	404	PEG	C4-C3-O2-C2
3	D	402	8YV	O2-C2-C3-O3
7	A	407	PG4	C3-C4-O3-C5
5	A	404	PEG	O2-C3-C4-O4
3	C	402	8YV	O4-C4-C5-C6
3	C	402	8YV	C3-C4-C5-C6
3	C	402	8YV	O4-C4-C5-O5
3	A	402	8YV	C2-C3-C4-C5
2	D	401	NAP	O4B-C4B-C5B-O5B
2	A	401	NAP	O4B-C4B-C5B-O5B
2	C	401	NAP	O4B-C4B-C5B-O5B
2	B	401	NAP	O4B-C4B-C5B-O5B

There are no ring outliers.

5 monomers are involved in 5 short contacts:

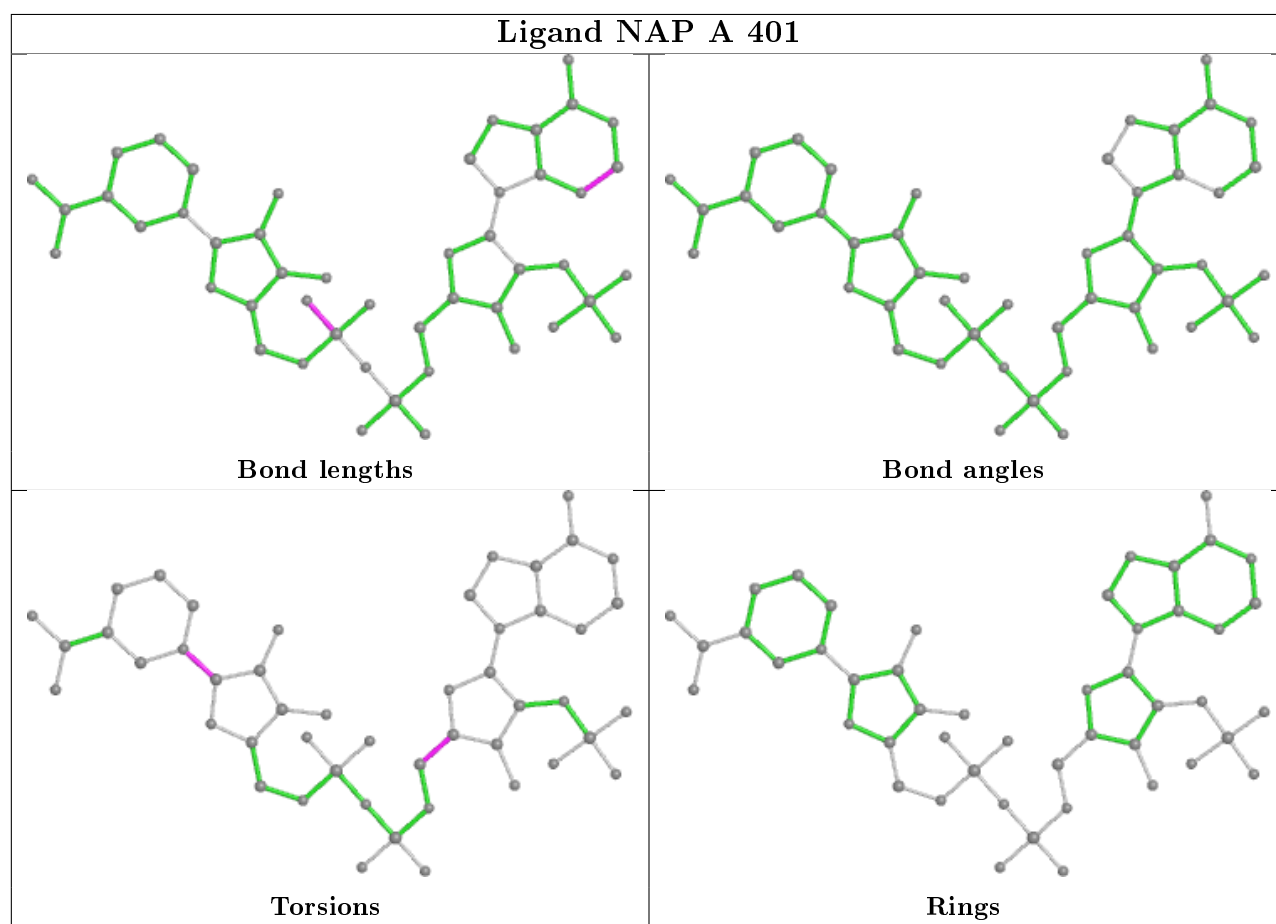
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	NAP	1	0

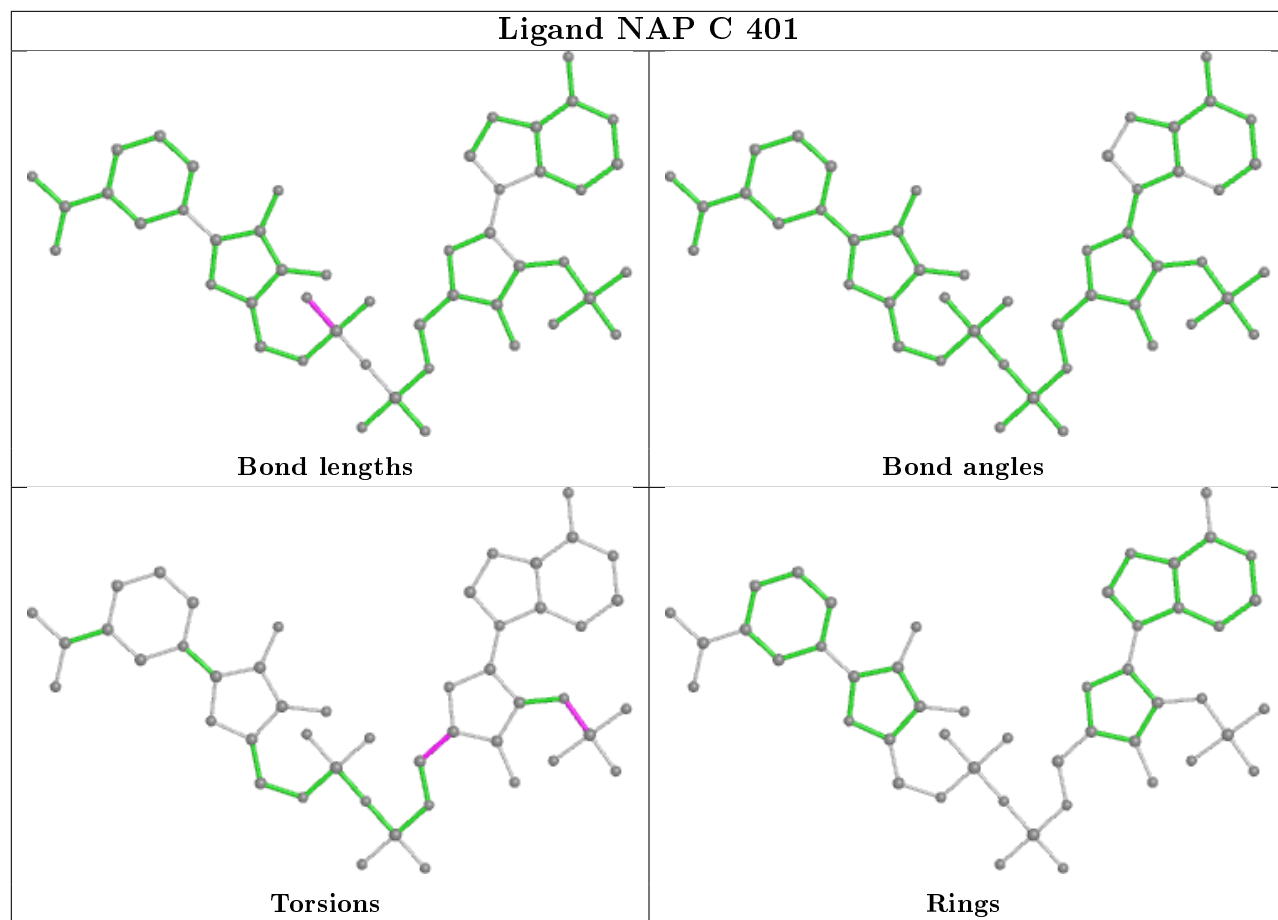
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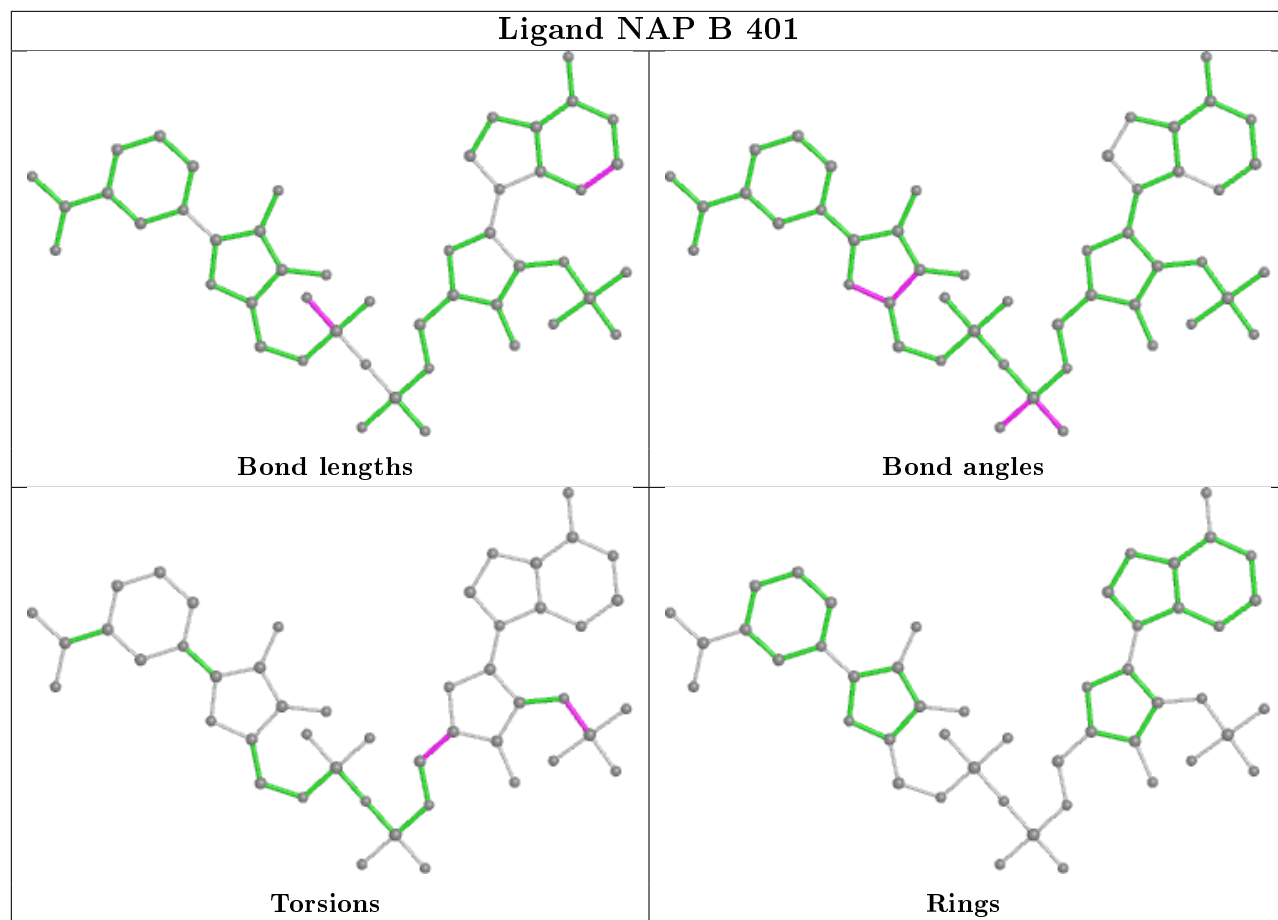
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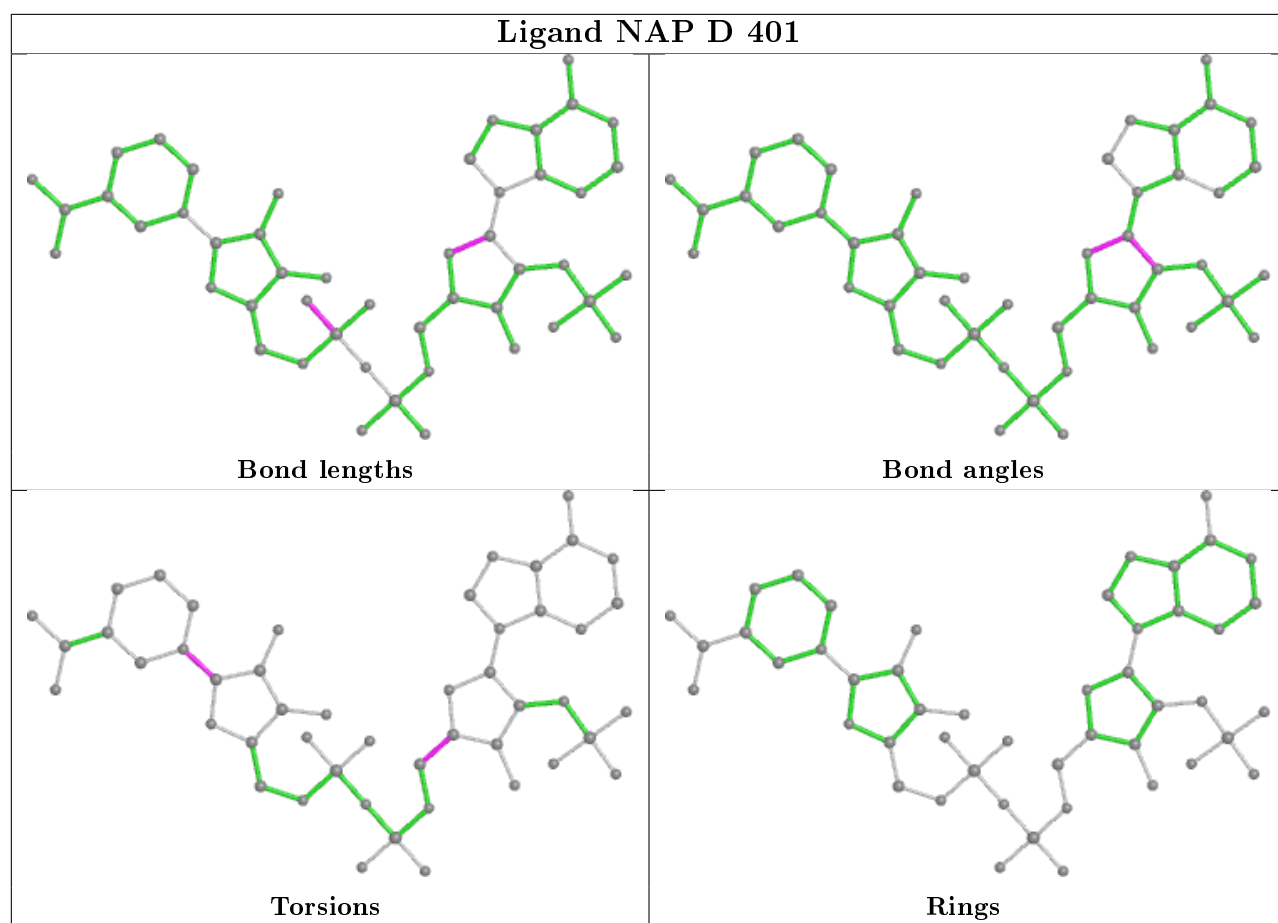
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	8YV	1	0
3	A	402	8YV	1	0
3	C	402	8YV	1	0
3	D	402	8YV	1	0

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	319/323 (98%)	-0.29	3 (0%) 84 89	16, 25, 42, 77	0
1	B	318/323 (98%)	-0.30	2 (0%) 89 92	18, 26, 41, 66	0
1	C	318/323 (98%)	-0.13	2 (0%) 89 92	22, 33, 50, 74	0
1	D	318/323 (98%)	-0.13	4 (1%) 77 83	23, 34, 51, 92	0
All	All	1273/1292 (98%)	-0.21	11 (0%) 84 89	16, 30, 48, 92	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	320	ILE	4.6
1	A	178	GLY	3.6
1	B	320	ILE	3.0
1	C	179	LEU	2.9
1	D	180	GLY	2.5
1	A	180	GLY	2.4
1	B	178	GLY	2.4
1	D	39	ALA	2.3
1	D	179	LEU	2.2
1	A	179	LEU	2.1
1	D	320	ILE	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 monosaccharides in this entry.

6.4 Ligands ⓘ

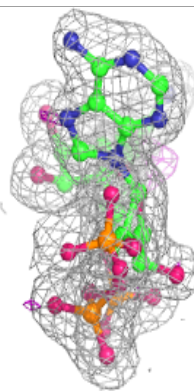
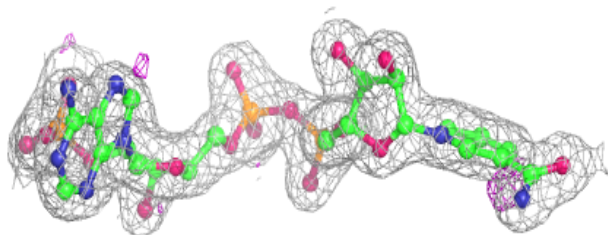
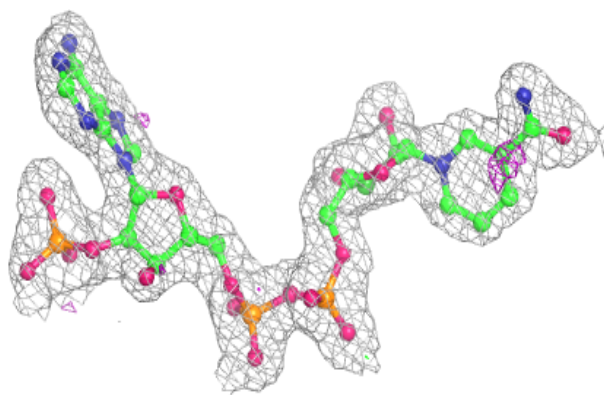
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	MG	A	408	1/1	0.71	0.11	56,56,56,56	0
3	8YV	D	402	13/13	0.82	0.13	30,38,46,50	0
5	PEG	D	404	7/7	0.82	0.15	53,55,63,66	0
8	MG	D	409	1/1	0.83	0.16	68,68,68,68	0
8	MG	B	408	1/1	0.83	0.14	55,55,55,55	0
8	MG	A	409	1/1	0.85	0.22	50,50,50,50	0
4	NA	D	403	1/1	0.87	0.10	37,37,37,37	0
8	MG	C	405	1/1	0.88	0.18	50,50,50,50	0
3	8YV	C	402	13/13	0.88	0.11	26,35,67,73	0
5	PEG	A	404	7/7	0.88	0.11	36,39,46,58	0
4	NA	A	403	1/1	0.89	0.08	31,31,31,31	0
8	MG	B	407	1/1	0.90	0.32	44,44,44,44	0
7	PG4	A	407	13/13	0.90	0.13	32,45,69,76	0
3	8YV	A	402	13/13	0.91	0.10	20,26,35,36	0
9	GOL	D	407	6/6	0.92	0.14	33,36,46,51	0
3	8YV	B	402	13/13	0.92	0.11	21,28,37,49	0
4	NA	C	403	1/1	0.92	0.06	46,46,46,46	0
8	MG	D	408	1/1	0.93	0.06	60,60,60,60	0
9	GOL	B	406	6/6	0.94	0.13	25,33,57,63	0
4	NA	B	403	1/1	0.95	0.07	29,29,29,29	0
2	NAP	C	401	48/48	0.96	0.08	22,30,38,42	0
6	CL	D	406	1/1	0.96	0.09	51,51,51,51	0
2	NAP	D	401	48/48	0.96	0.08	20,29,41,43	0
6	CL	A	406	1/1	0.97	0.07	41,41,41,41	0
2	NAP	A	401	48/48	0.97	0.07	17,23,33,37	0
2	NAP	B	401	48/48	0.97	0.07	15,22,30,33	0
6	CL	B	405	1/1	0.99	0.07	41,41,41,41	0
6	CL	B	404	1/1	0.99	0.06	26,26,26,26	0
6	CL	A	405	1/1	0.99	0.05	27,27,27,27	0
6	CL	D	405	1/1	0.99	0.06	35,35,35,35	0
6	CL	C	404	1/1	0.99	0.06	34,34,34,34	0

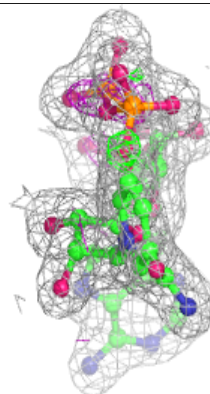
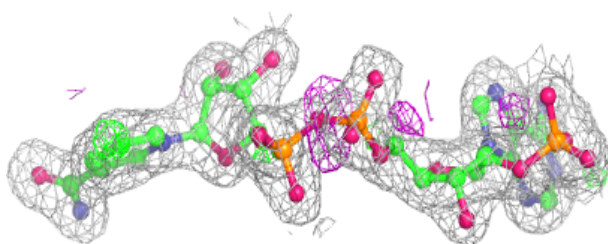
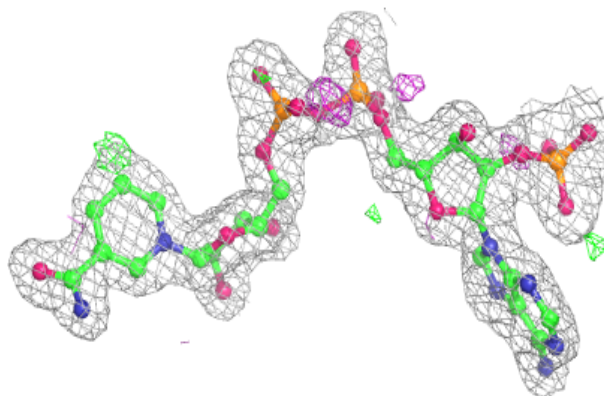
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAP C 401:

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

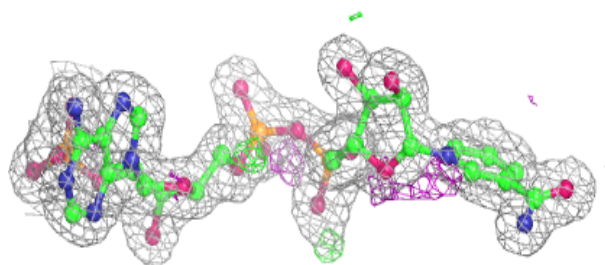
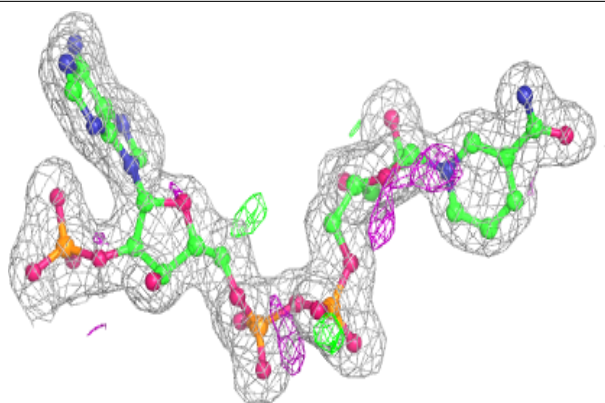
**Electron density around NAP D 401:**

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

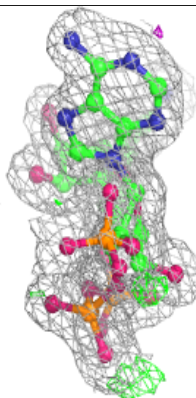
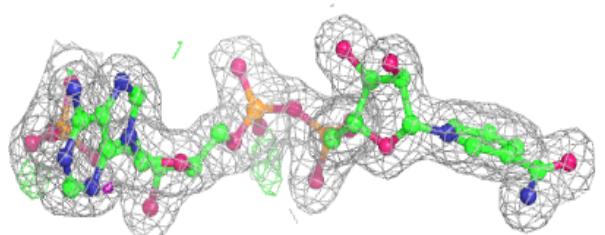
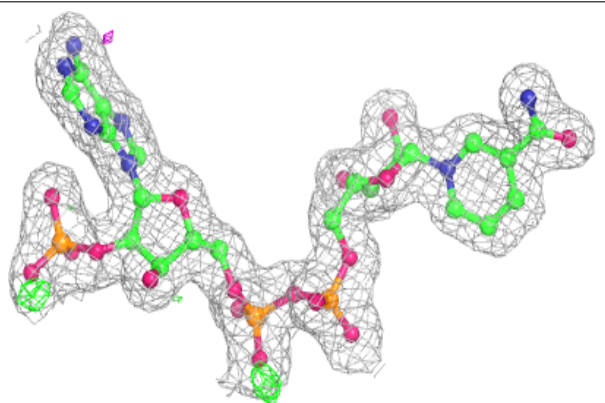


Electron density around NAP A 401:

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

**Electron density around NAP B 401:**

$2mF_o-DF_c$ (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.