



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 01:08 PM UTC

PDB ID : 3WID / pdb\_00003wid  
Title : Structure of a glucose dehydrogenase T277F mutant in complex with NADP  
Authors : Sakuraba, H.; Kanoh, Y.; Yoneda, K.; Ohshima, T.  
Deposited on : 2013-09-10  
Resolution : 2.25 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

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<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>.

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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

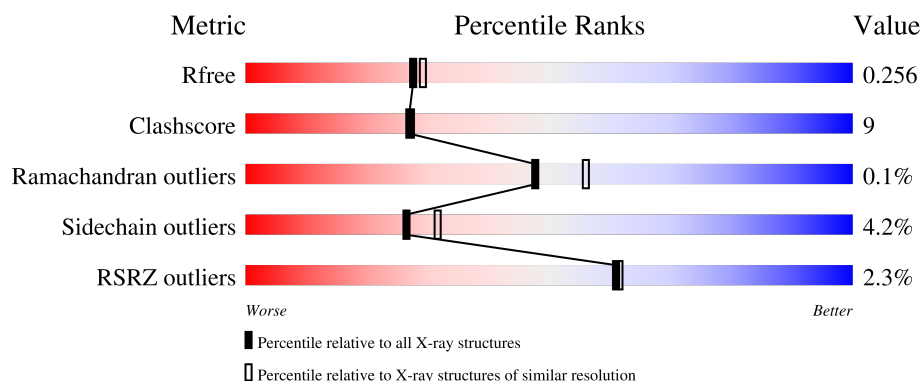
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*




The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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  $\geq 3$ , 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	369	 80% 16% ..
1	B	369	 81% 14% ..
1	C	369	 4% 80% 16% ..
1	D	369	 5% 77% 18% ..

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
5	1PE	A	1004	-	-	X	-
5	1PE	B	4005	-	-	X	-
5	1PE	B	4006	-	-	X	-
5	1PE	C	1004	-	-	X	-
5	1PE	C	1005	-	-	X	-
5	1PE	D	4001	-	-	X	-
5	1PE	D	4005	-	-	X	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11962 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 Glucose 1-dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	360	Total	C	N	O	S	0	0	0
			2827	1804	473	537	13			
1	B	360	Total	C	N	O	S	0	0	0
			2827	1804	473	537	13			
1	C	358	Total	C	N	O	S	0	0	0
			2809	1793	470	533	13			
1	D	360	Total	C	N	O	S	0	0	0
			2827	1804	473	537	13			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	277	PHE	THR	engineered mutation	UNP Q979W2
A	362	LEU	-	expression tag	UNP Q979W2
A	363	GLU	-	expression tag	UNP Q979W2
A	364	HIS	-	expression tag	UNP Q979W2
A	365	HIS	-	expression tag	UNP Q979W2
A	366	HIS	-	expression tag	UNP Q979W2
A	367	HIS	-	expression tag	UNP Q979W2
A	368	HIS	-	expression tag	UNP Q979W2
A	369	HIS	-	expression tag	UNP Q979W2
B	277	PHE	THR	engineered mutation	UNP Q979W2
B	362	LEU	-	expression tag	UNP Q979W2
B	363	GLU	-	expression tag	UNP Q979W2
B	364	HIS	-	expression tag	UNP Q979W2
B	365	HIS	-	expression tag	UNP Q979W2
B	366	HIS	-	expression tag	UNP Q979W2
B	367	HIS	-	expression tag	UNP Q979W2
B	368	HIS	-	expression tag	UNP Q979W2
B	369	HIS	-	expression tag	UNP Q979W2
C	277	PHE	THR	engineered mutation	UNP Q979W2
C	362	LEU	-	expression tag	UNP Q979W2
C	363	GLU	-	expression tag	UNP Q979W2

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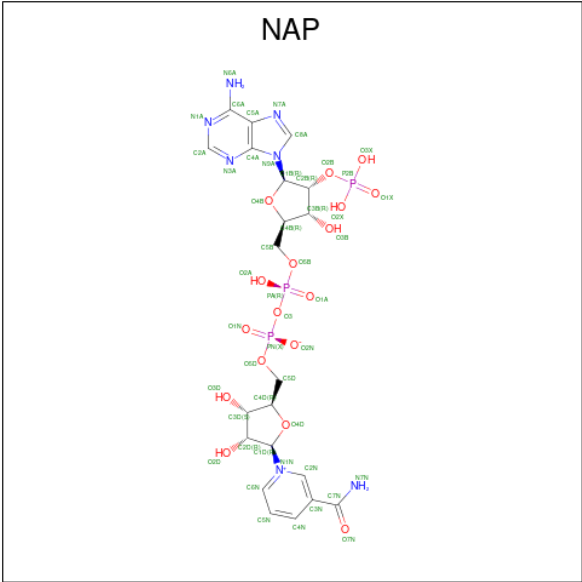
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Chain	Residue	Modelled	Actual	Comment	Reference
C	364	HIS	-	expression tag	UNP Q979W2
C	365	HIS	-	expression tag	UNP Q979W2
C	366	HIS	-	expression tag	UNP Q979W2
C	367	HIS	-	expression tag	UNP Q979W2
C	368	HIS	-	expression tag	UNP Q979W2
C	369	HIS	-	expression tag	UNP Q979W2
D	277	PHE	THR	engineered mutation	UNP Q979W2
D	362	LEU	-	expression tag	UNP Q979W2
D	363	GLU	-	expression tag	UNP Q979W2
D	364	HIS	-	expression tag	UNP Q979W2
D	365	HIS	-	expression tag	UNP Q979W2
D	366	HIS	-	expression tag	UNP Q979W2
D	367	HIS	-	expression tag	UNP Q979W2
D	368	HIS	-	expression tag	UNP Q979W2
D	369	HIS	-	expression tag	UNP Q979W2

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

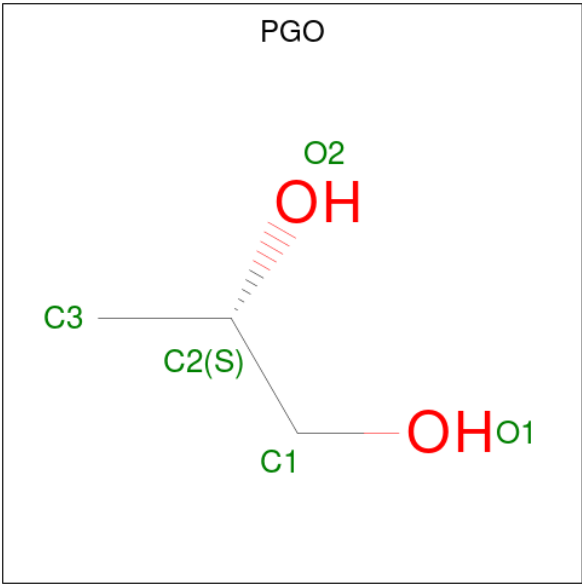
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: C<sub>21</sub>H<sub>28</sub>N<sub>7</sub>O<sub>17</sub>P<sub>3</sub>).



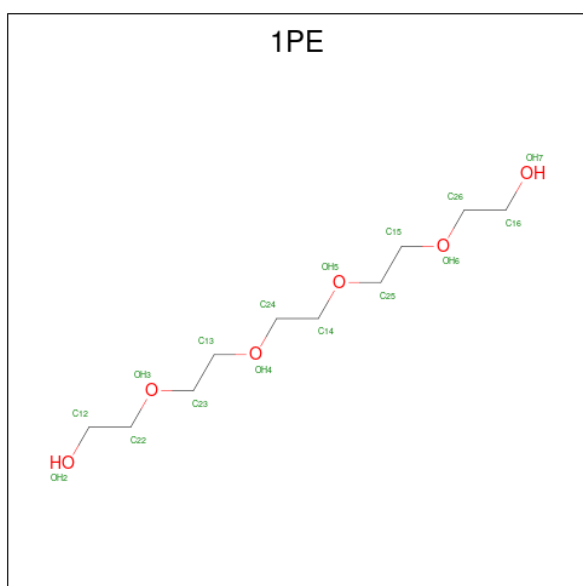
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 4 is S-1,2-PROPANEDIOL (CCD ID: PGO) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			5	3	2		
4	B	1	Total	C	O	0	0
			5	3	2		
4	C	1	Total	C	O	0	0
			5	3	2		
4	D	1	Total	C	O	0	0
			5	3	2		

- Molecule 5 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula:  $C_{10}H_{22}O_6$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			16	10	6		
5	B	1	Total	C	O	0	0
			16	10	6		
5	B	1	Total	C	O	0	0
			16	10	6		
5	B	1	Total	C	O	0	0
			16	10	6		
5	C	1	Total	C	O	0	0
			16	10	6		
5	C	1	Total	C	O	0	0
			16	10	6		
5	D	1	Total	C	O	0	0
			16	10	6		
5	D	1	Total	C	O	0	0
			16	10	6		

- Molecule 6 is water.

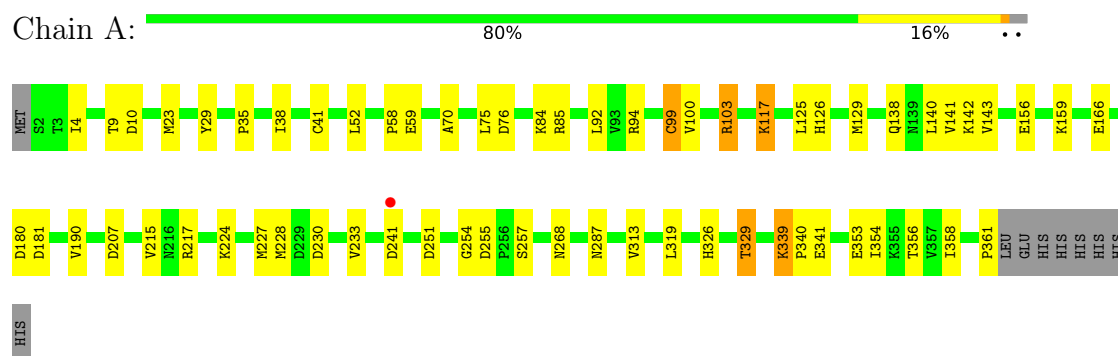
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	92	Total 92	O 92	0	0
6	B	109	Total 109	O 109	0	0
6	C	58	Total 58	O 58	0	0
6	D	69	Total 69	O 69	0	0



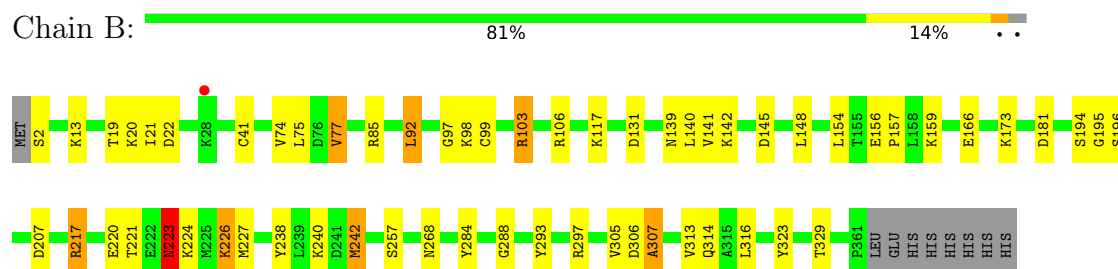
### 3 Residue-property plots [i](#)

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.

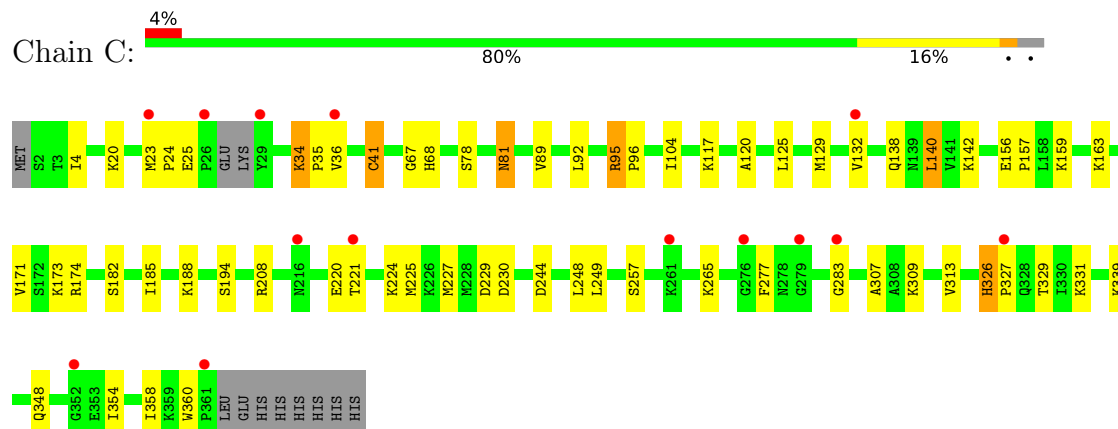
#### • Molecule 1: Glucose 1-dehydrogenase



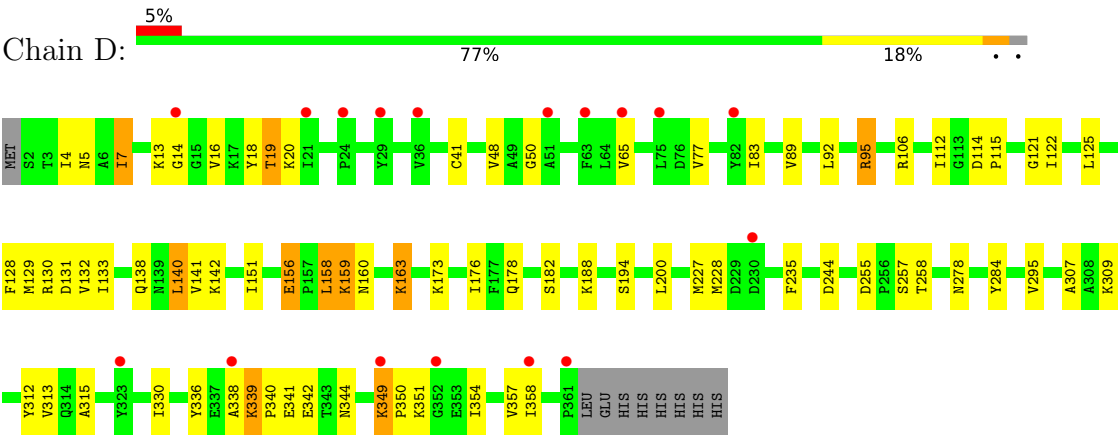
#### • Molecule 1: Glucose 1-dehydrogenase



#### • Molecule 1: Glucose 1-dehydrogenase



● Molecule 1: Glucose 1-dehydrogenase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.17Å 90.35Å 120.37Å 90.00° 90.84° 90.00°	Depositor
Resolution (Å)	39.27 – 2.25 39.27 – 2.25	Depositor EDS
% Data completeness (in resolution range)	94.3 (39.27-2.25) 94.3 (39.27-2.25)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	8.60 (at 2.24Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, $R_{free}$	0.197 , 0.251 (Not available) , 0.256	Depositor DCC
$R_{free}$ test set	3881 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	39.1	Xtriage
Anisotropy	0.478	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 46.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.012 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	11962	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.82% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PGO, ZN, NAP, 1PE

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	1.21	5/2887 (0.2%)	1.10	5/3904 (0.1%)
1	B	1.29	7/2887 (0.2%)	1.14	3/3904 (0.1%)
1	C	1.09	4/2868 (0.1%)	1.11	5/3878 (0.1%)
1	D	1.07	2/2887 (0.1%)	1.12	5/3904 (0.1%)
All	All	1.17	18/11529 (0.2%)	1.12	18/15590 (0.1%)

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	70	ALA	CA-C	-6.81	1.44	1.52
1	A	207	ASP	C-O	-6.39	1.16	1.24
1	B	207	ASP	C-O	-6.23	1.16	1.24
1	B	145	ASP	C-O	-5.85	1.17	1.24
1	A	319	LEU	C-O	-5.83	1.17	1.24

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	14	GLY	N-CA-C	7.39	119.84	110.38
1	B	240	LYS	N-CA-C	7.38	119.40	111.36
1	D	156	GLU	CA-C-N	-7.16	111.60	119.19
1	D	156	GLU	C-N-CA	-7.16	111.60	119.19
1	A	255	ASP	CA-C-N	6.66	126.17	119.24

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	2827	0	2808	42	0
1	B	2827	0	2808	50	0
1	C	2809	0	2788	46	0
1	D	2827	0	2808	65	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	48	0	25	1	0
3	B	48	0	25	2	0
3	C	48	0	25	1	0
3	D	48	0	25	0	0
4	A	5	0	8	3	0
4	B	5	0	8	0	0
4	C	5	0	8	0	0
4	D	5	0	8	3	0
5	A	16	0	22	9	0
5	B	48	0	66	24	0
5	C	32	0	44	19	0
5	D	32	0	44	18	0
6	A	92	0	0	3	0
6	B	109	0	0	5	0
6	C	58	0	0	8	0
6	D	69	0	0	3	0
All	All	11962	0	11520	212	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 9.

The worst 5 of 212 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:106:ARG:HH11	5:B:4005:1PE:H241	1.18	1.03
1:D:227:MET:HE1	1:D:354:ILE:CD1	1.93	0.98
1:B:106:ARG:HD3	5:B:4005:1PE:H241	1.49	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:227:MET:HE1	1:D:354:ILE:HD11	1.48	0.91
1:A:100:VAL:HA	1:A:103:ARG:HD3	1.49	0.91

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	358/369 (97%)	347 (97%)	10 (3%)	1 (0%)	36	40
1	B	358/369 (97%)	351 (98%)	7 (2%)	0	100	100
1	C	354/369 (96%)	336 (95%)	18 (5%)	0	100	100
1	D	358/369 (97%)	339 (95%)	19 (5%)	0	100	100
All	All	1428/1476 (97%)	1373 (96%)	54 (4%)	1 (0%)	48	56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	CYS

### 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	313/322 (97%)	302 (96%)	11 (4%)	32	39

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	313/322 (97%)	302 (96%)	11 (4%)	32	39
1	C	311/322 (97%)	293 (94%)	18 (6%)	18	19
1	D	313/322 (97%)	300 (96%)	13 (4%)	26	31
All	All	1250/1288 (97%)	1197 (96%)	53 (4%)	26	31

5 of 53 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	92	LEU
1	C	249	LEU
1	D	257	SER
1	C	140	LEU
1	C	225	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 23 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	326	HIS
1	D	5	ASN
1	C	348	GLN
1	D	118	HIS
1	B	178	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry

Of 20 ligands modelled in this entry, 4 are monoatomic - leaving 16 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
5	1PE	C	1005	-	15,15,15	1.08	0	14,14,14	1.21	2 (14%)
4	PGO	D	4004	-	4,4,4	0.42	0	4,4,4	1.21	1 (25%)
3	NAP	A	1002	-	50,52,52	1.36	6 (12%)	71,80,80	1.72	12 (16%)
3	NAP	C	1002	-	50,52,52	1.29	5 (10%)	71,80,80	1.87	24 (33%)
3	NAP	B	4003	-	50,52,52	1.34	6 (12%)	71,80,80	2.20	24 (33%)
4	PGO	A	1003	-	4,4,4	0.48	0	4,4,4	1.13	1 (25%)
5	1PE	A	1004	-	15,15,15	0.90	0	14,14,14	1.16	1 (7%)
5	1PE	C	1004	-	15,15,15	0.74	0	14,14,14	0.74	0
5	1PE	B	4006	-	15,15,15	1.08	1 (6%)	14,14,14	0.98	0
3	NAP	D	4003	-	50,52,52	1.39	6 (12%)	71,80,80	1.78	12 (16%)
5	1PE	D	4001	-	15,15,15	0.87	0	14,14,14	0.59	0
5	1PE	B	4005	-	15,15,15	1.07	0	14,14,14	1.07	0
5	1PE	D	4005	-	15,15,15	0.72	0	14,14,14	0.89	1 (7%)
4	PGO	C	1003	-	4,4,4	0.67	0	4,4,4	0.75	0
5	1PE	B	4001	-	15,15,15	1.09	1 (6%)	14,14,14	0.99	0
4	PGO	B	4004	-	4,4,4	0.48	0	4,4,4	0.69	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
5	1PE	C	1005	-	-	7/13/13/13	-
4	PGO	D	4004	-	-	2/2/2/2	-
3	NAP	A	1002	-	-	10/35/67/67	0/5/5/5
3	NAP	C	1002	-	-	6/35/67/67	0/5/5/5
3	NAP	B	4003	-	-	9/35/67/67	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PGO	A	1003	-	-	0/2/2/2	-
5	1PE	A	1004	-	-	4/13/13/13	-
5	1PE	C	1004	-	-	9/13/13/13	-
5	1PE	B	4006	-	-	6/13/13/13	-
3	NAP	D	4003	-	-	8/35/67/67	0/5/5/5
5	1PE	D	4001	-	-	8/13/13/13	-
5	1PE	B	4005	-	-	6/13/13/13	-
5	1PE	D	4005	-	-	6/13/13/13	-
4	PGO	C	1003	-	-	0/2/2/2	-
5	1PE	B	4001	-	-	6/13/13/13	-
4	PGO	B	4004	-	-	0/2/2/2	-

The worst 5 of 25 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	4003	NAP	C8A-N9A	-3.73	1.31	1.37
3	D	4003	NAP	C5A-N7A	-3.69	1.32	1.39
3	B	4003	NAP	PA-O3	3.68	1.63	1.59
3	A	1002	NAP	C4A-N9A	-3.59	1.30	1.37
3	B	4003	NAP	C5A-C4A	3.30	1.45	1.39

The worst 5 of 78 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4003	NAP	C5A-C4A-N3A	-6.78	117.38	126.72
3	B	4003	NAP	C4A-N9A-C8A	5.97	112.01	105.74
3	B	4003	NAP	O7N-C7N-C3N	5.73	126.60	119.60
3	D	4003	NAP	N3A-C4A-N9A	5.49	136.50	127.17
3	A	1002	NAP	C5A-C4A-N3A	-5.46	119.19	126.72

There are no chirality outliers.

5 of 87 torsion outliers are listed below:

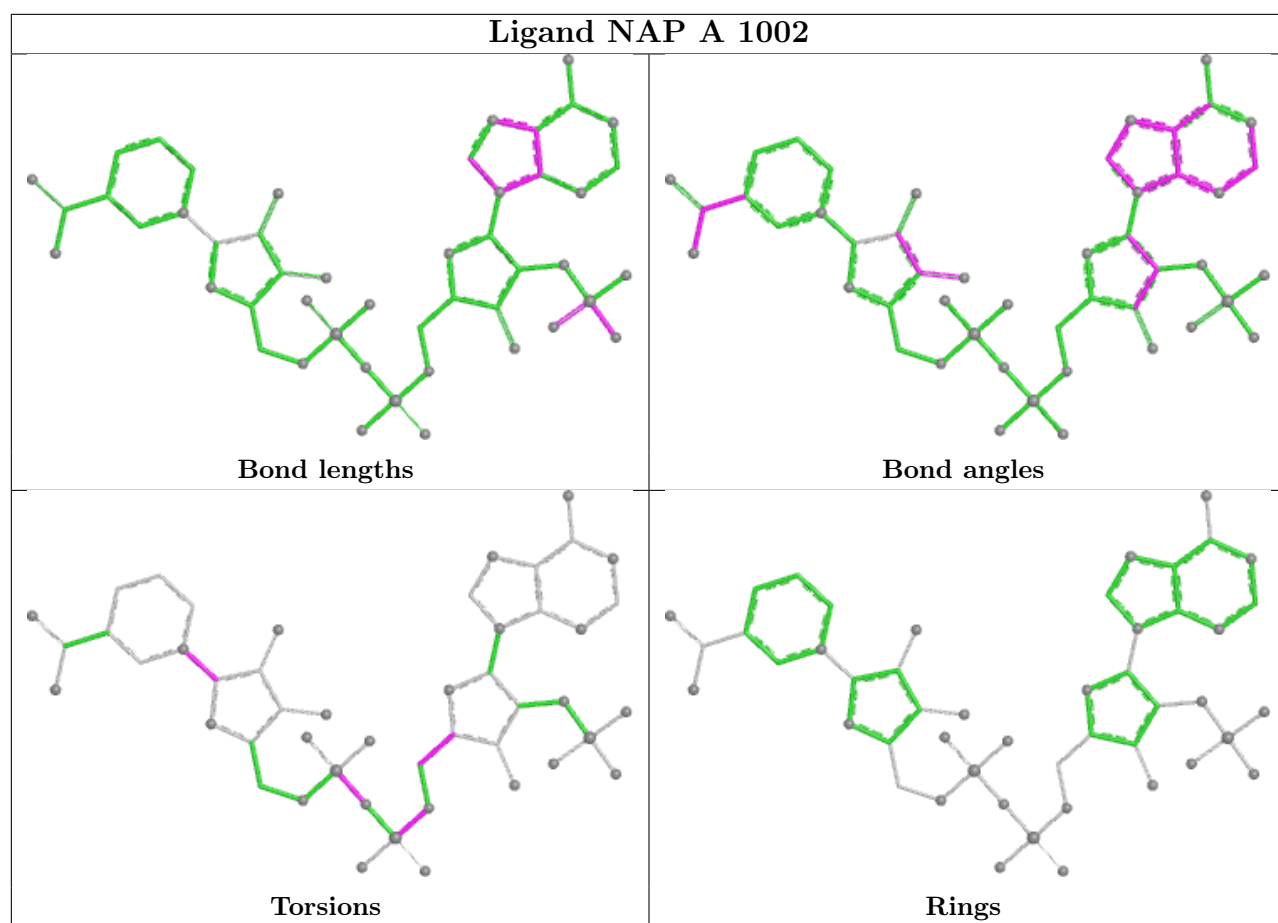
Mol	Chain	Res	Type	Atoms
3	A	1002	NAP	C5B-O5B-PA-O2A
3	A	1002	NAP	C5B-O5B-PA-O3
3	A	1002	NAP	O4D-C1D-N1N-C2N
3	A	1002	NAP	O4D-C1D-N1N-C6N
3	A	1002	NAP	C2D-C1D-N1N-C2N

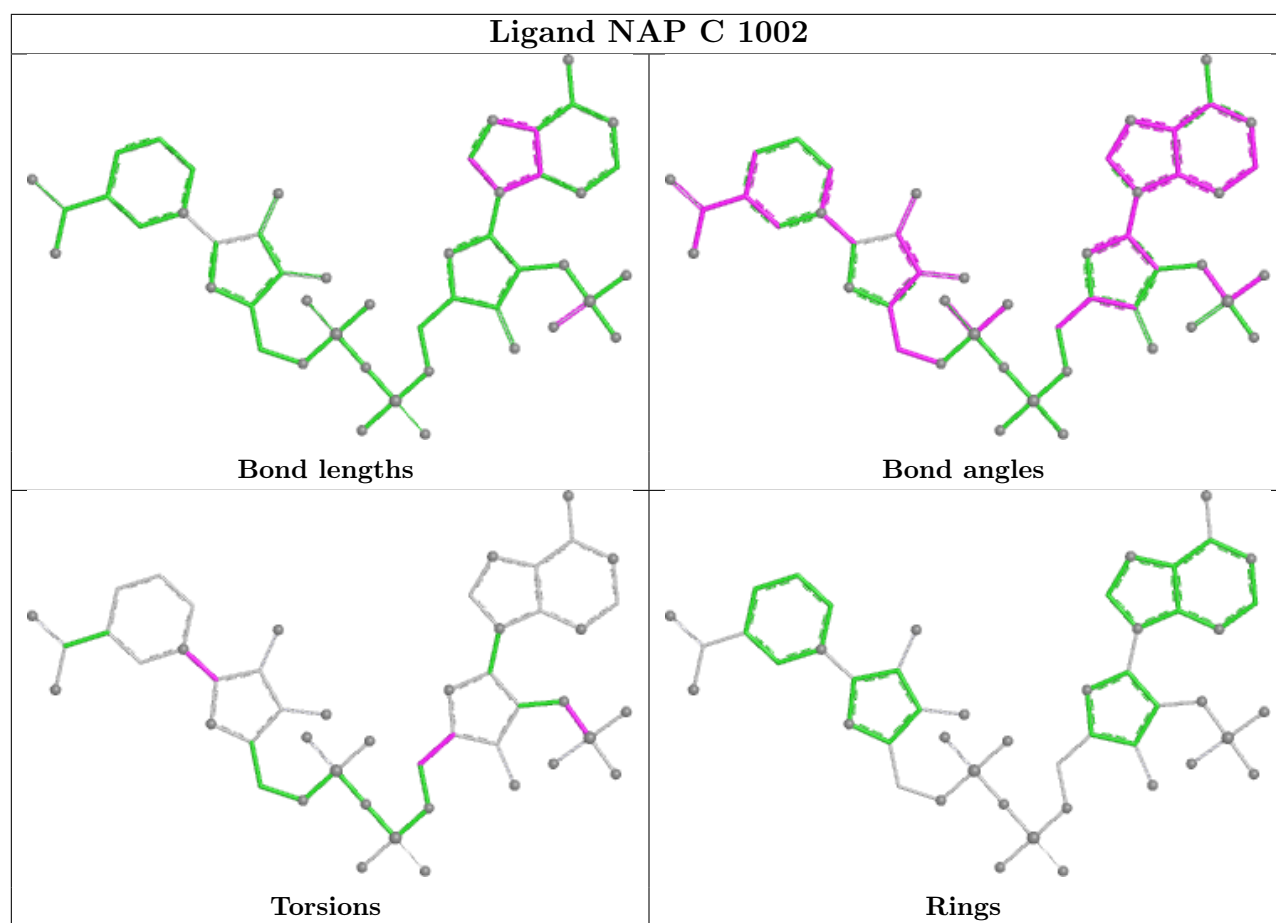
There are no ring outliers.

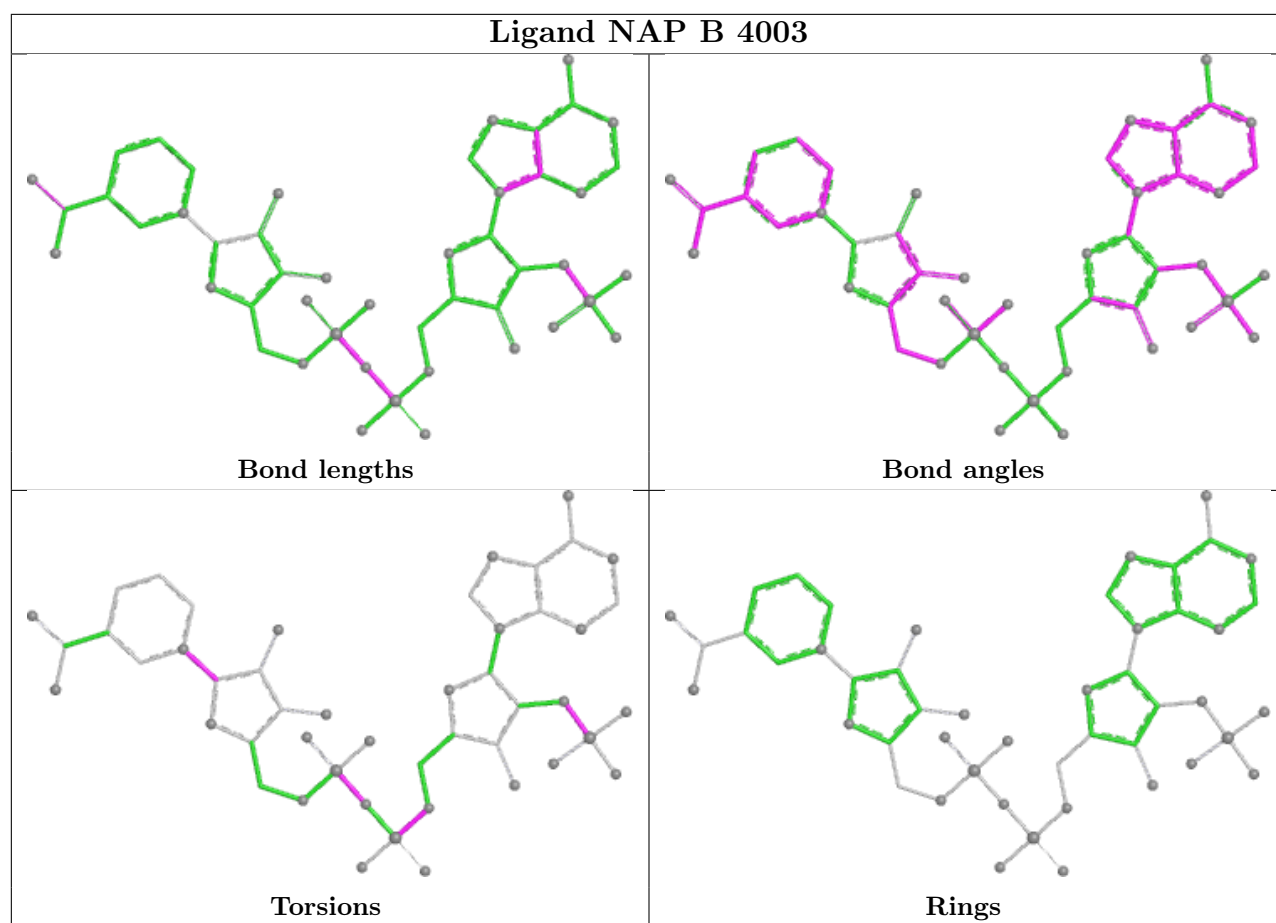
13 monomers are involved in 80 short contacts:

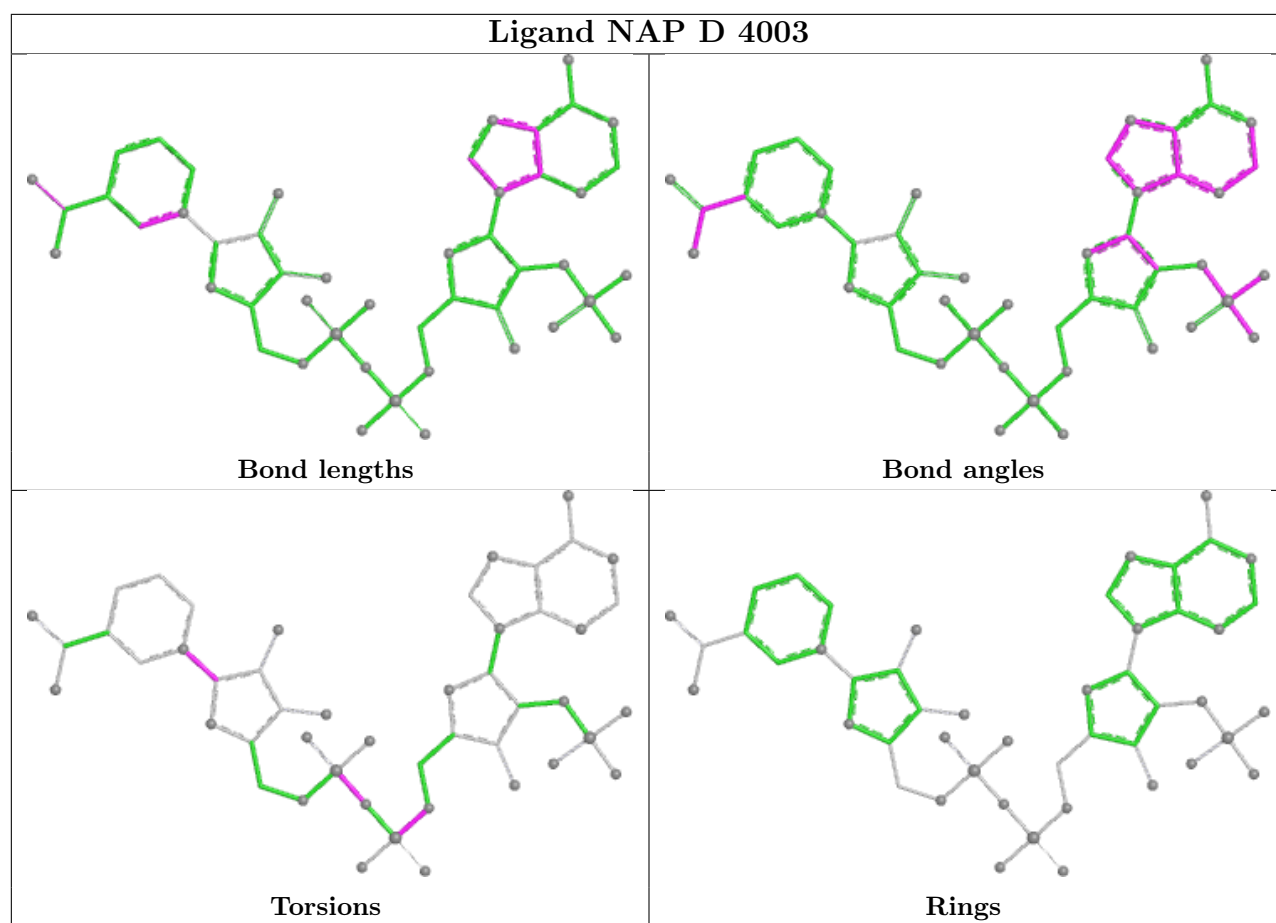
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	1005	1PE	11	0
4	D	4004	PGO	3	0
3	A	1002	NAP	1	0
3	C	1002	NAP	1	0
3	B	4003	NAP	2	0
4	A	1003	PGO	3	0
5	A	1004	1PE	9	0
5	C	1004	1PE	8	0
5	B	4006	1PE	9	0
5	D	4001	1PE	7	0
5	B	4005	1PE	12	0
5	D	4005	1PE	11	0
5	B	4001	1PE	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	360/369 (97%)	0.09	1 (0%) 90 90	27, 43, 64, 82	0
1	B	360/369 (97%)	-0.10	1 (0%) 90 90	24, 39, 61, 93	0
1	C	358/369 (97%)	0.43	14 (3%) 43 43	31, 50, 78, 104	0
1	D	360/369 (97%)	0.52	17 (4%) 36 35	26, 52, 85, 99	0
All	All	1438/1476 (97%)	0.23	33 (2%) 61 61	24, 45, 76, 104	0

The worst 5 of 33 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	352	GLY	3.9
1	D	14	GLY	3.8
1	C	279	GLY	3.6
1	C	26	PRO	3.2
1	C	352	GLY	2.9

### 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, 95<sup>th</sup> 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.

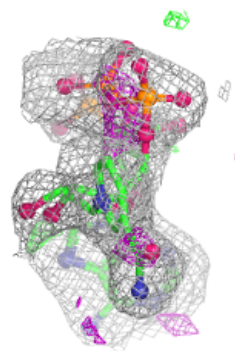
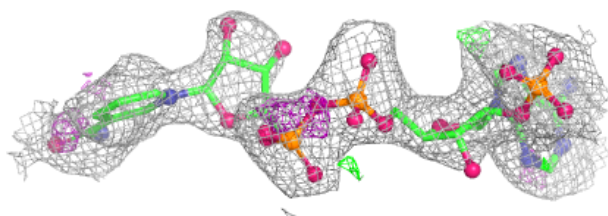
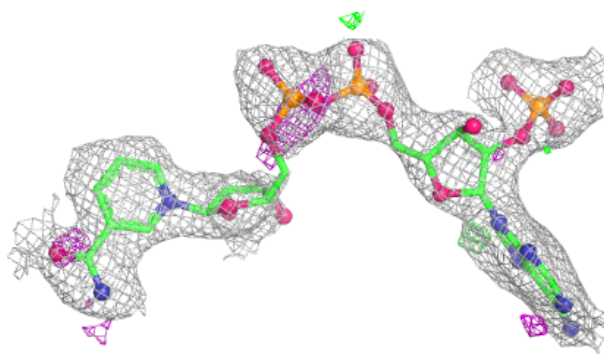
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	1PE	B	4005	16/16	0.75	0.17	43,56,63,64	0
5	1PE	D	4005	16/16	0.75	0.16	46,54,69,72	0
4	PGO	D	4004	5/5	0.80	0.12	48,52,57,61	0
5	1PE	B	4006	16/16	0.81	0.13	38,47,53,55	0
5	1PE	B	4001	16/16	0.82	0.13	44,52,56,57	0
4	PGO	C	1003	5/5	0.82	0.14	42,48,54,59	0
5	1PE	D	4001	16/16	0.83	0.14	37,51,61,62	0
5	1PE	A	1004	16/16	0.83	0.13	42,50,61,64	0
5	1PE	C	1005	16/16	0.85	0.14	34,50,62,62	0
5	1PE	C	1004	16/16	0.87	0.11	41,55,63,65	0
3	NAP	C	1002	48/48	0.88	0.13	43,62,75,78	0
3	NAP	D	4003	48/48	0.89	0.12	37,54,66,68	0
3	NAP	A	1002	48/48	0.90	0.11	42,55,61,66	0
3	NAP	B	4003	48/48	0.93	0.10	39,50,68,71	0
4	PGO	A	1003	5/5	0.94	0.08	45,47,49,51	0
4	PGO	B	4004	5/5	0.96	0.10	43,43,47,51	0
2	ZN	A	1001	1/1	0.99	0.02	36,36,36,36	0
2	ZN	B	4002	1/1	0.99	0.02	31,31,31,31	0
2	ZN	C	1001	1/1	0.99	0.02	34,34,34,34	0
2	ZN	D	4002	1/1	0.99	0.02	38,38,38,38	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.

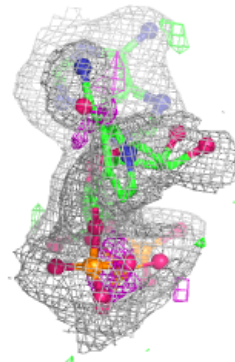
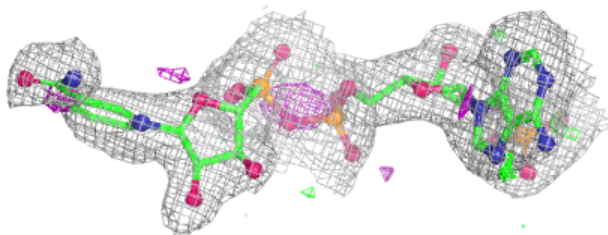
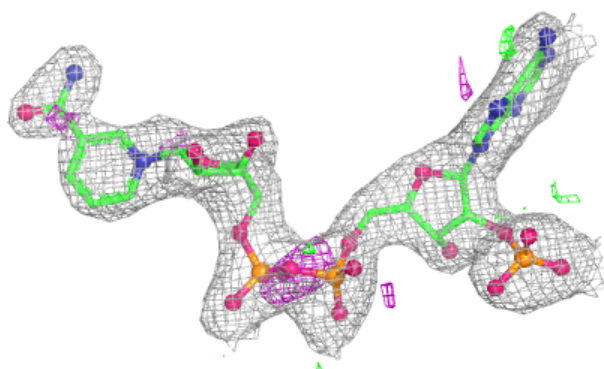


**Electron density around NAP C 1002:**

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

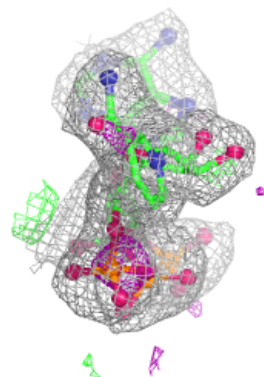
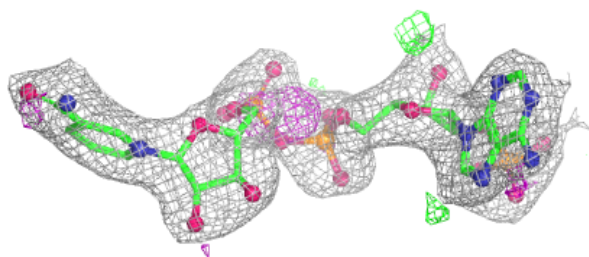
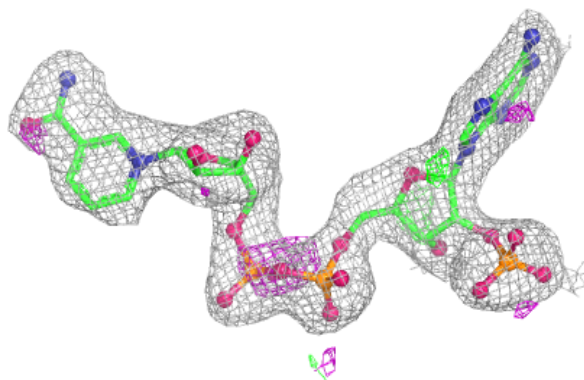
**Electron density around NAP D 4003:**

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

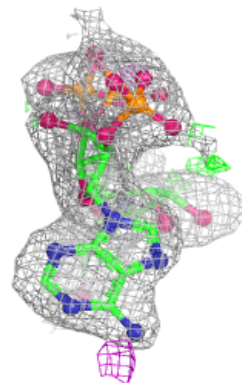
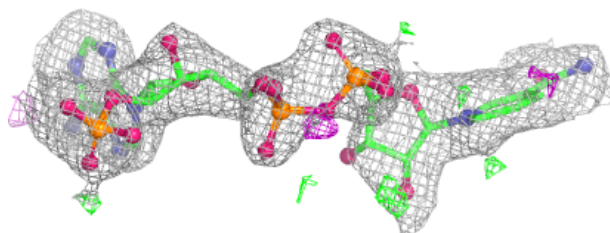
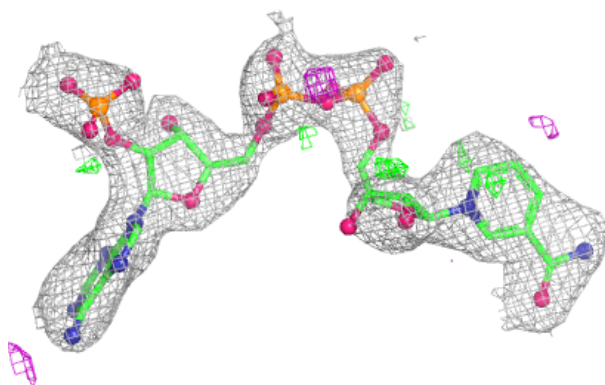


**Electron density around NAP A 1002:**

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

**Electron density around NAP B 4003:**

$2mF_o-DF_c$  (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.