



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 16, 2023 – 11:12 AM EDT

PDB ID : 2GRA
Title : crystal structure of Human Pyrroline-5-carboxylate Reductase complexed with nadp
Authors : Meng, Z.; Lou, Z.; Liu, Z.; Rao, Z.
Deposited on : 2006-04-23
Resolution : 3.10 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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

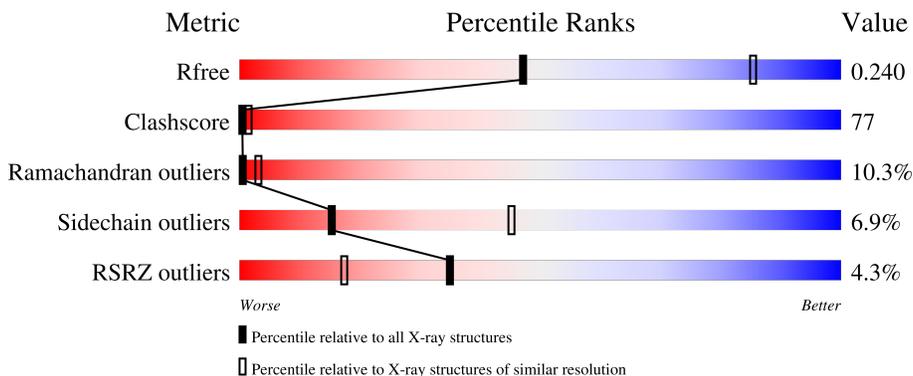
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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	277	 11% 25% 58% 14%
1	B	277	 11% 21% 64% 15%
1	C	277	 10% 27% 64% 9%
1	D	277	 29% 60% 12%
1	E	277	 25% 64% 11%

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	NAP	A	1300	X	-	-	-
2	NAP	B	2300	-	-	-	X
2	NAP	C	3300	X	-	-	X
2	NAP	D	4300	-	-	-	X
2	NAP	E	5300	X	-	-	X
3	GLU	B	2301	-	-	-	X
3	GLU	C	3301	-	-	-	X
3	GLU	D	4301	-	-	-	X
3	GLU	E	5301	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11079 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.

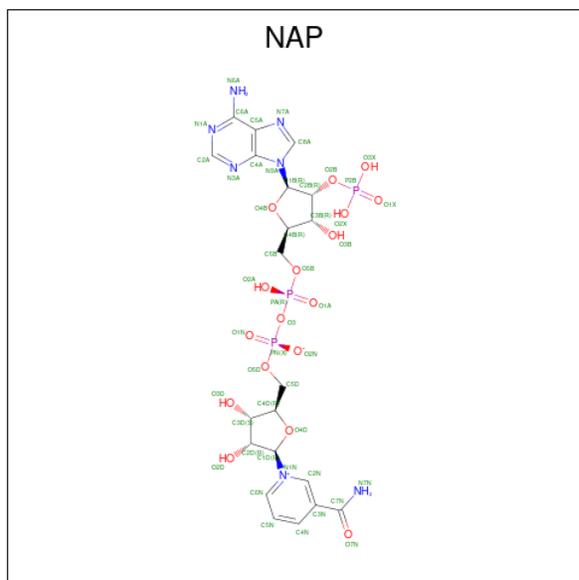
- Molecule 1 is a protein called Pyrroline-5-carboxylate reductase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	277	2038	1279	363	383	13	0	0	0
1	B	276	2023	1270	358	382	13	0	0	0
1	C	277	2032	1276	360	383	13	0	0	0
1	D	277	2038	1279	363	383	13	0	0	0
1	E	277	2038	1279	363	383	13	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

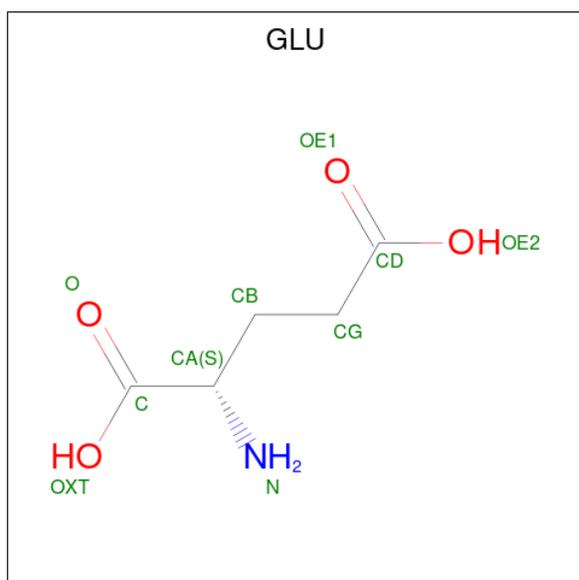
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ARG	-	cloning artifact	UNP P32322
A	0	GLY	-	cloning artifact	UNP P32322
B	-1	ARG	-	cloning artifact	UNP P32322
B	0	GLY	-	cloning artifact	UNP P32322
C	-1	ARG	-	cloning artifact	UNP P32322
C	0	GLY	-	cloning artifact	UNP P32322
D	-1	ARG	-	cloning artifact	UNP P32322
D	0	GLY	-	cloning artifact	UNP P32322
E	-1	ARG	-	cloning artifact	UNP P32322
E	0	GLY	-	cloning artifact	UNP P32322

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



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

- Molecule 3 is GLUTAMIC ACID (three-letter code: GLU) (formula: C₅H₉NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			10	5	1	4		
3	B	1	Total	C	N	O	0	0
			10	5	1	4		
3	C	1	Total	C	N	O	0	0
			10	5	1	4		
3	D	1	Total	C	N	O	0	0
			10	5	1	4		
3	E	1	Total	C	N	O	0	0
			10	5	1	4		

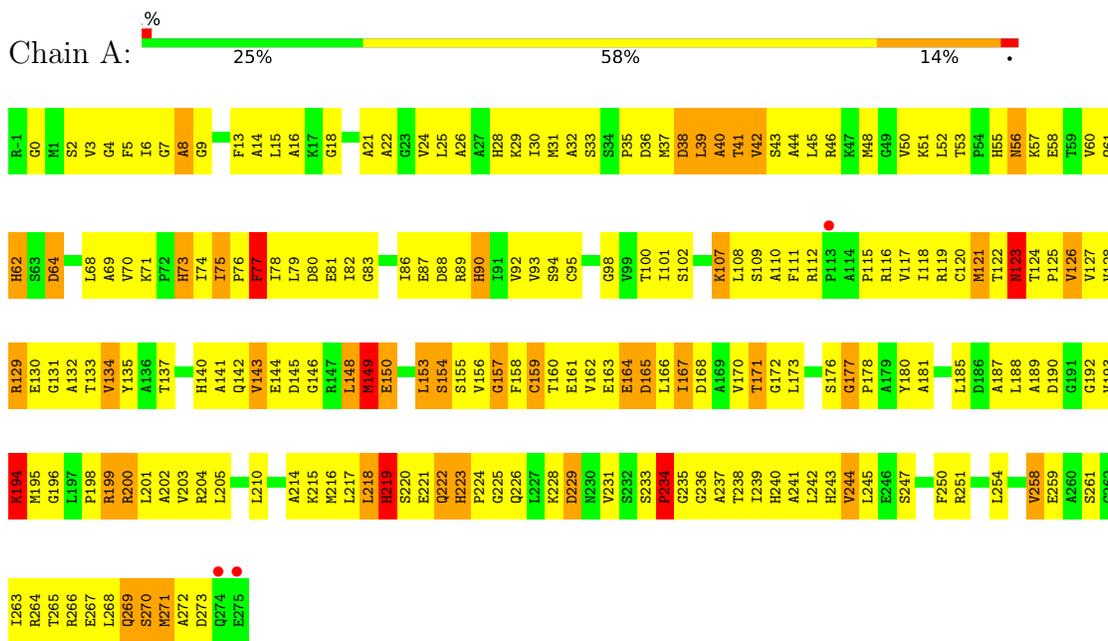
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	127	Total	O	0	0
			127	127		
4	B	100	Total	O	0	0
			100	100		
4	C	104	Total	O	0	0
			104	104		
4	D	125	Total	O	0	0
			125	125		
4	E	164	Total	O	0	0
			164	164		

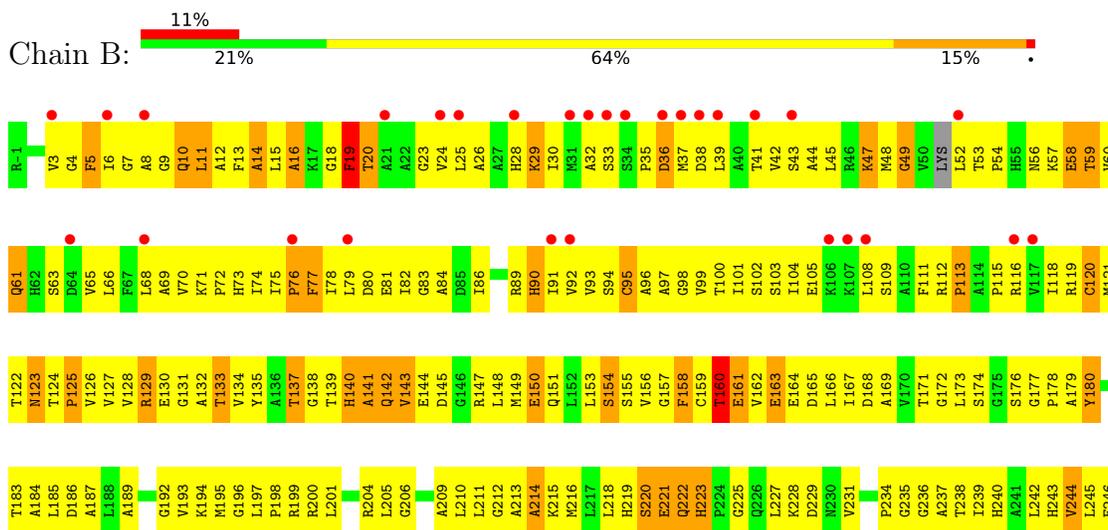
3 Residue-property plots

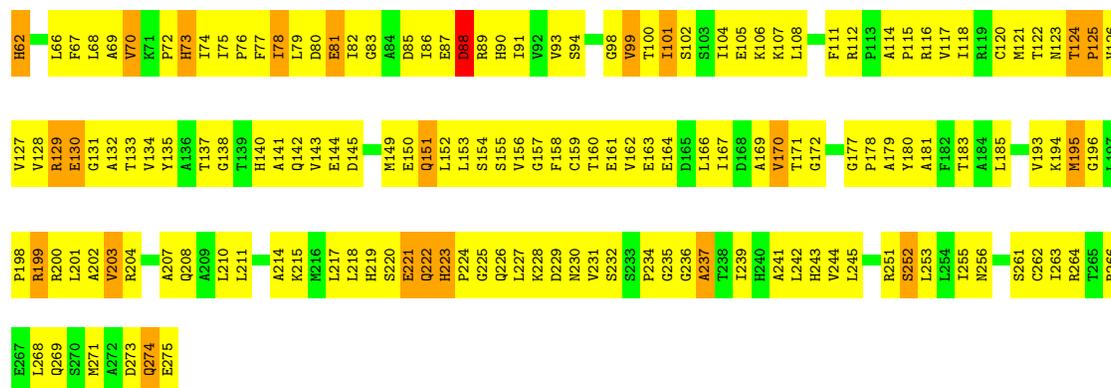
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Pyrroline-5-carboxylate reductase 1



- Molecule 1: Pyrroline-5-carboxylate reductase 1





4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	207.87Å 123.50Å 120.83Å 90.00° 121.97° 90.00°	Depositor
Resolution (Å)	50.00 – 3.10 48.42 – 3.01	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-3.10) 96.0 (48.42-3.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 3.01Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.228 , 0.249 0.223 , 0.240	Depositor DCC
R_{free} test set	2528 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	58.0	Xtrriage
Anisotropy	0.425	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 126.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	0.017 for $-1/2^*h+1/2^*k+1, 1/2^*h-1/2^*k+1, 1/2^*h+1/2^*k$ 0.026 for $-1/2^*h-1/2^*k+1, -1/2^*h-1/2^*k-1, 1/2^*h-1/2^*k$	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	11079	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.74% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.61	3/2069 (0.1%)	0.79	7/2800 (0.2%)
1	B	0.34	0/2053	0.65	1/2779 (0.0%)
1	C	0.35	0/2063	0.66	0/2793
1	D	0.36	0/2069	0.68	0/2800
1	E	0.39	0/2069	0.70	0/2800
All	All	0.42	3/10323 (0.0%)	0.70	8/13972 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	219	HIS	C-N	16.69	1.72	1.34
1	A	157	GLY	C-N	9.66	1.56	1.34
1	A	153	LEU	C-N	-9.30	1.12	1.34

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	219	HIS	O-C-N	-10.31	106.20	122.70
1	A	219	HIS	CA-C-N	7.21	133.06	117.20
1	A	157	GLY	O-C-N	7.14	134.12	122.70
1	A	219	HIS	C-N-CA	6.86	138.84	121.70
1	A	154	SER	O-C-N	6.26	132.72	122.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	218	LEU	Mainchain

5.2 Too-close contacts [i](#)

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	2038	0	2082	327	0
1	B	2023	0	2059	345	0
1	C	2032	0	2073	310	2
1	D	2038	0	2084	279	0
1	E	2038	0	2084	328	1
2	A	48	0	23	18	0
2	B	48	0	22	15	0
2	C	48	0	21	12	0
2	D	48	0	22	11	1
2	E	48	0	21	14	0
3	A	10	0	5	3	0
3	B	10	0	5	3	0
3	C	10	0	5	1	0
3	D	10	0	5	2	0
3	E	10	0	5	0	0
4	A	127	0	0	63	3
4	B	100	0	0	45	0
4	C	104	0	0	57	1
4	D	125	0	0	59	0
4	E	164	0	0	53	0
All	All	11079	0	10516	1619	4

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 77.

The worst 5 of 1619 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)
2:E:5300:NAP:N7A	2:E:5300:NAP:C8A	1.72	1.51
2:A:1300:NAP:N7A	2:A:1300:NAP:C8A	1.72	1.49
2:C:3300:NAP:C8A	2:C:3300:NAP:N7A	1.72	1.45
2:D:4300:NAP:C8A	2:D:4300:NAP:N7A	1.71	1.40
2:B:2300:NAP:N7A	2:B:2300:NAP:C8A	1.72	1.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:1340:HOH:O	4:C:3362:HOH:O[2_555]	1.63	0.57
1:C:129:ARG:CG	4:A:1320:HOH:O[2_555]	2.05	0.15
1:E:208:GLN:OE1	2:D:4300:NAP:O2N[2_555]	2.16	0.04
1:C:129:ARG:CD	4:A:1320:HOH:O[2_555]	2.19	0.01

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	275/277 (99%)	188 (68%)	62 (22%)	25 (9%)	1	4
1	B	272/277 (98%)	167 (61%)	67 (25%)	38 (14%)	0	1
1	C	275/277 (99%)	193 (70%)	59 (22%)	23 (8%)	1	5
1	D	275/277 (99%)	193 (70%)	51 (18%)	31 (11%)	0	2
1	E	275/277 (99%)	185 (67%)	66 (24%)	24 (9%)	1	4
All	All	1372/1385 (99%)	926 (68%)	305 (22%)	141 (10%)	0	3

5 of 141 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8	ALA
1	A	39	LEU

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Mol	Chain	Res	Type
1	A	107	LYS
1	A	129	ARG
1	A	143	VAL

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	215/215 (100%)	190 (88%)	25 (12%)	5 22
1	B	213/215 (99%)	198 (93%)	15 (7%)	15 45
1	C	214/215 (100%)	204 (95%)	10 (5%)	26 59
1	D	215/215 (100%)	204 (95%)	11 (5%)	24 56
1	E	215/215 (100%)	202 (94%)	13 (6%)	19 49
All	All	1072/1075 (100%)	998 (93%)	74 (7%)	15 45

5 of 74 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	144	GLU
1	E	195	MET
1	D	211	LEU
1	E	62	HIS
1	A	271	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 26 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	240	HIS
1	D	55	HIS
1	E	219	HIS
1	D	28	HIS
1	D	90	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

10 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	NAP	E	5300	-	45,52,52	4.75	23 (51%)	56,80,80	3.29	22 (39%)
3	GLU	A	1301	-	8,9,9	0.97	0	10,11,11	0.92	0
3	GLU	C	3301	-	8,9,9	0.84	0	10,11,11	0.84	0
3	GLU	D	4301	-	8,9,9	1.02	0	10,11,11	0.92	0
3	GLU	E	5301	-	8,9,9	0.99	0	10,11,11	0.91	0
2	NAP	D	4300	1	45,52,52	4.79	24 (53%)	56,80,80	3.15	23 (41%)
2	NAP	B	2300	-	45,52,52	4.95	25 (55%)	56,80,80	3.16	23 (41%)
3	GLU	B	2301	-	8,9,9	1.01	0	10,11,11	1.00	0
2	NAP	C	3300	-	45,52,52	4.88	24 (53%)	56,80,80	3.64	22 (39%)
2	NAP	A	1300	-	45,52,52	5.33	25 (55%)	56,80,80	4.16	31 (55%)

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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAP	E	5300	-	1/1/12/12	6/31/67/67	0/5/5/5
3	GLU	A	1301	-	-	1/9/9/9	-
3	GLU	C	3301	-	-	0/9/9/9	-
3	GLU	D	4301	-	-	6/9/9/9	-
3	GLU	E	5301	-	-	2/9/9/9	-
2	NAP	D	4300	1	-	6/31/67/67	0/5/5/5
2	NAP	B	2300	-	-	6/31/67/67	0/5/5/5
3	GLU	B	2301	-	-	4/9/9/9	-
2	NAP	C	3300	-	1/1/12/12	6/31/67/67	0/5/5/5
2	NAP	A	1300	-	1/1/12/12	5/31/67/67	0/5/5/5

The worst 5 of 121 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2300	NAP	C8A-N7A	21.40	1.72	1.34
2	E	5300	NAP	C8A-N7A	21.32	1.72	1.34
2	A	1300	NAP	C8A-N7A	21.10	1.72	1.34
2	C	3300	NAP	C8A-N7A	20.98	1.72	1.34
2	D	4300	NAP	C8A-N7A	20.86	1.71	1.34

The worst 5 of 121 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	3300	NAP	O5B-C5B-C4B	12.75	152.88	108.99
2	C	3300	NAP	C2N-N1N-C1D	-12.18	92.01	119.14
2	A	1300	NAP	O7N-C7N-N7N	-11.61	106.09	122.58
2	E	5300	NAP	C2N-N1N-C1D	-11.05	94.53	119.14
2	A	1300	NAP	O5B-C5B-C4B	10.13	143.86	108.99

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	1300	NAP	C1D
2	C	3300	NAP	C1D
2	E	5300	NAP	C1D

5 of 42 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1300	NAP	C2D-C1D-N1N-C2N
2	A	1300	NAP	C2D-C1D-N1N-C6N

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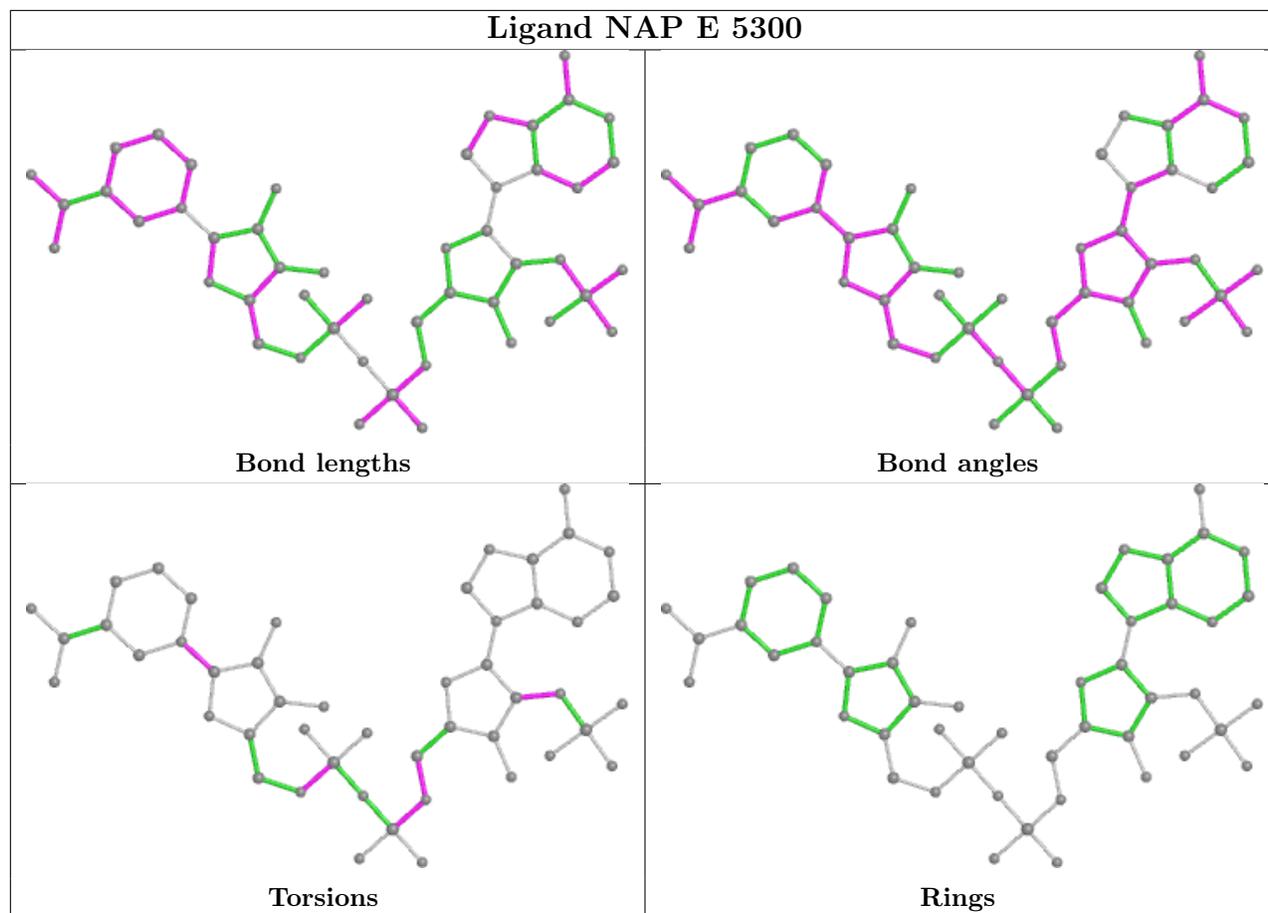
Mol	Chain	Res	Type	Atoms
2	B	2300	NAP	C2D-C1D-N1N-C2N
2	B	2300	NAP	C2D-C1D-N1N-C6N
2	C	3300	NAP	C4B-C5B-O5B-PA

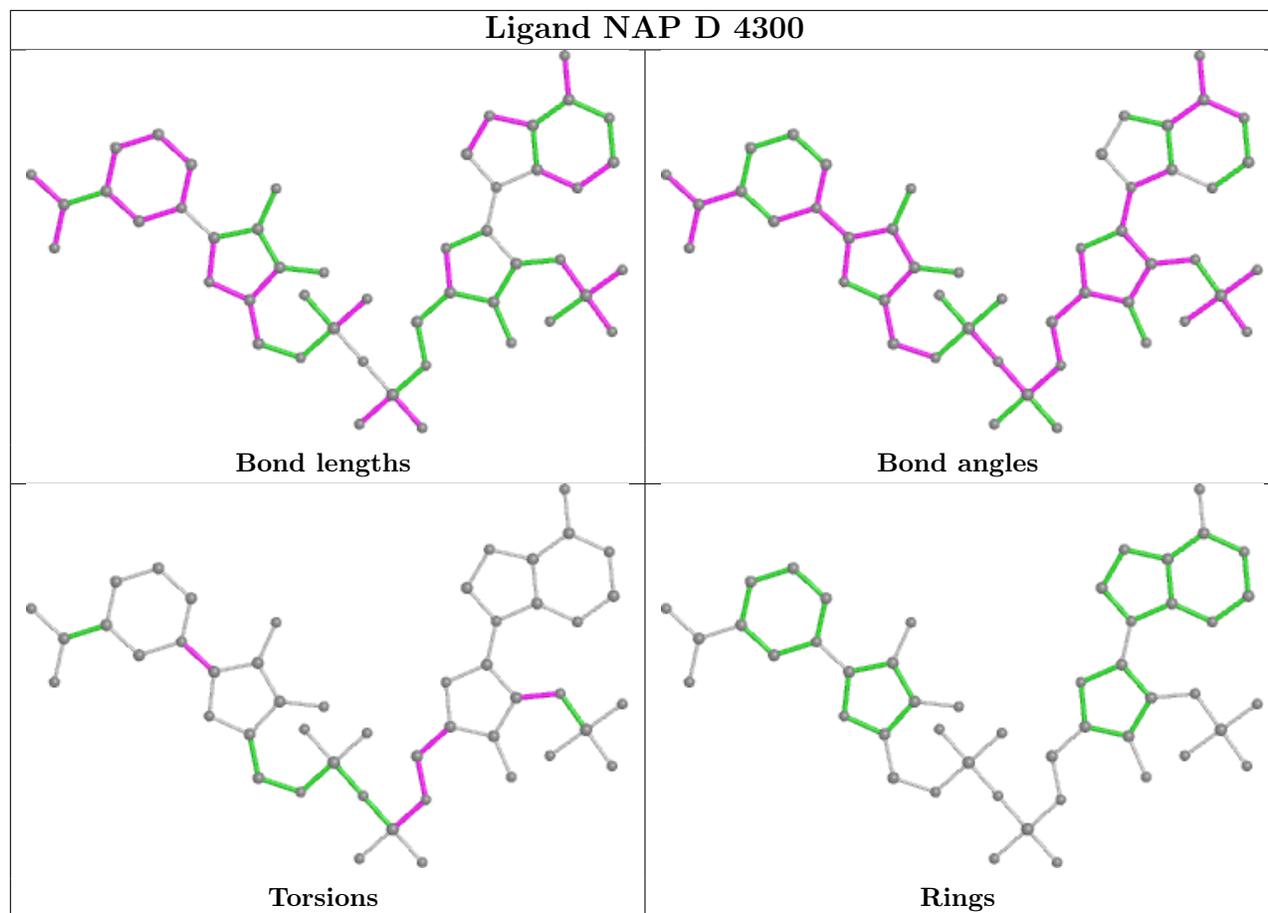
There are no ring outliers.

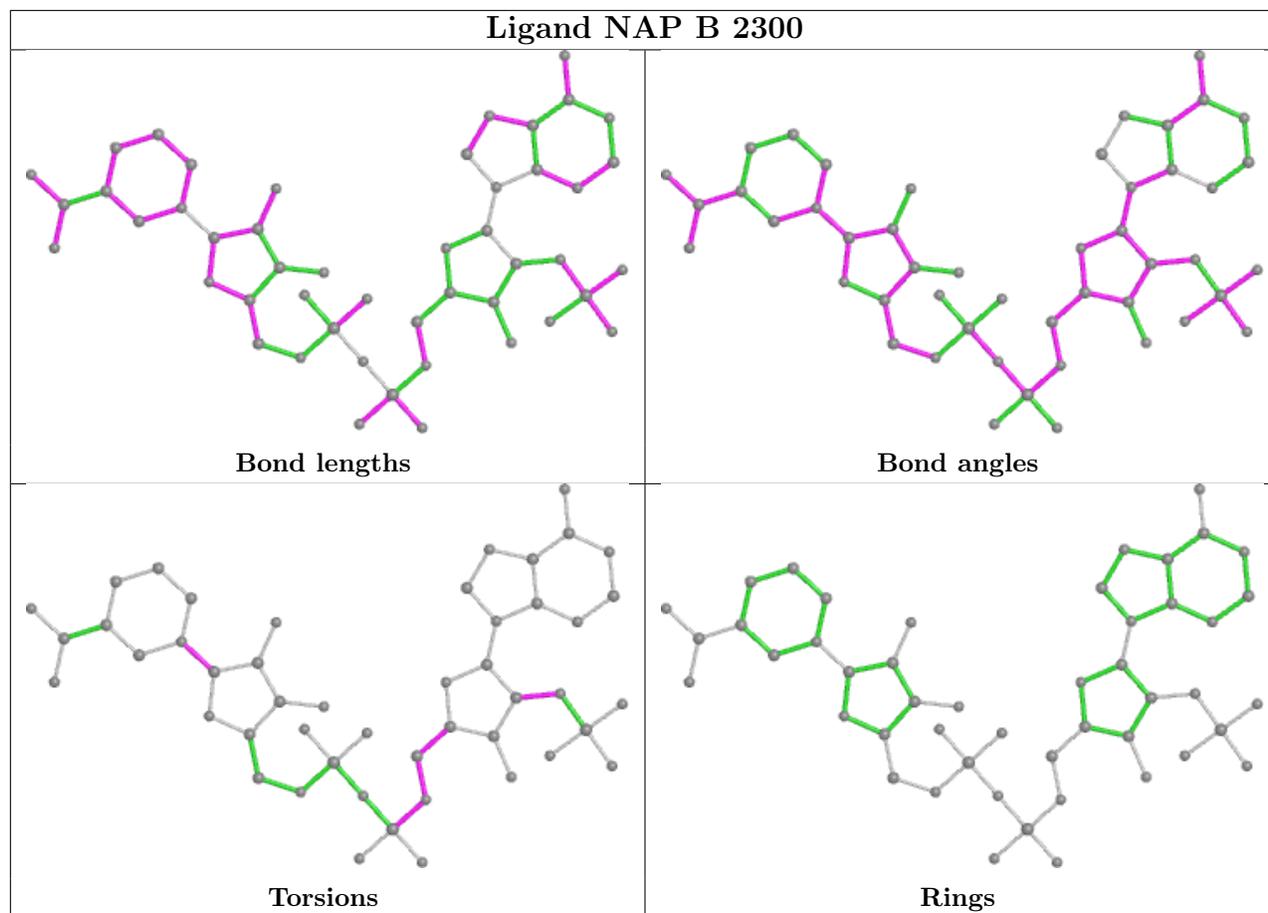
9 monomers are involved in 73 short contacts:

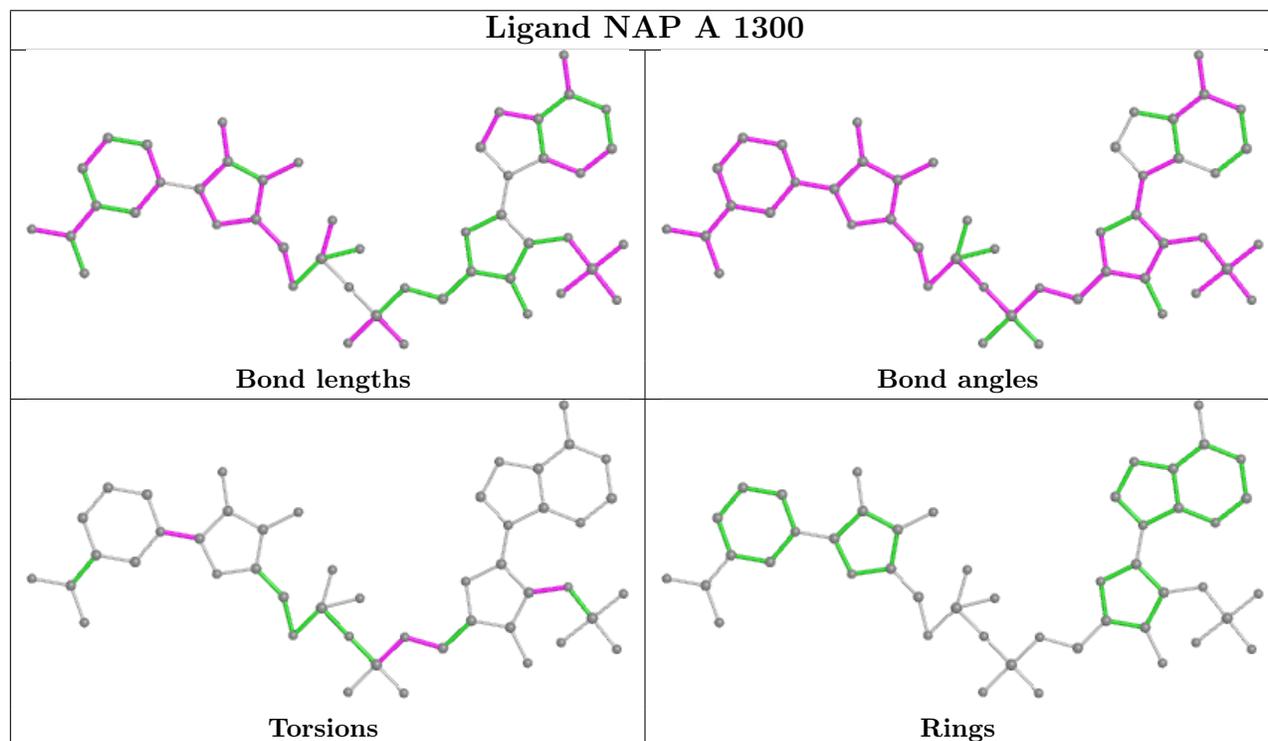
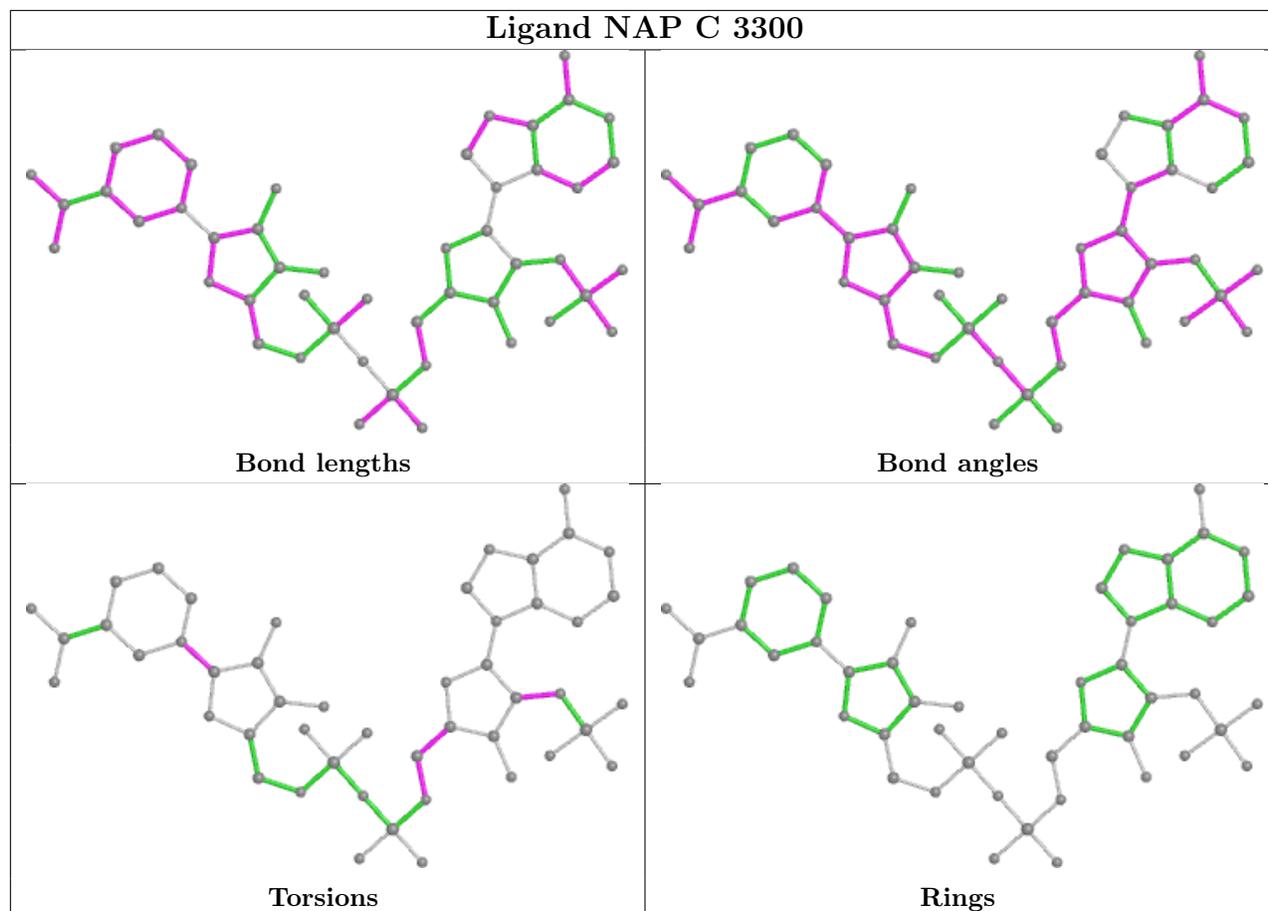
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	5300	NAP	14	0
3	A	1301	GLU	3	0
3	C	3301	GLU	1	0
3	D	4301	GLU	2	0
2	D	4300	NAP	11	1
2	B	2300	NAP	15	0
3	B	2301	GLU	3	0
2	C	3300	NAP	12	0
2	A	1300	NAP	18	0

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	219:HIS	C	220:SER	N	1.72
1	A	153:LEU	C	154:SER	N	1.12

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, 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	277/277 (100%)	-0.47	3 (1%) 80 64	20, 68, 115, 173	0
1	B	276/277 (99%)	0.28	30 (10%) 5 2	19, 115, 183, 200	0
1	C	277/277 (100%)	0.24	27 (9%) 7 2	16, 112, 178, 200	0
1	D	277/277 (100%)	-0.56	0 100 100	15, 62, 113, 178	0
1	E	277/277 (100%)	-0.53	0 100 100	14, 67, 124, 175	0
All	All	1384/1385 (99%)	-0.21	60 (4%) 35 17	14, 73, 170, 200	0

The worst 5 of 60 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	91	ILE	4.5
1	B	38	ASP	4.4
1	B	52	LEU	4.2
1	C	25	LEU	4.1
1	C	91	ILE	3.9

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 monosaccharides in this entry.

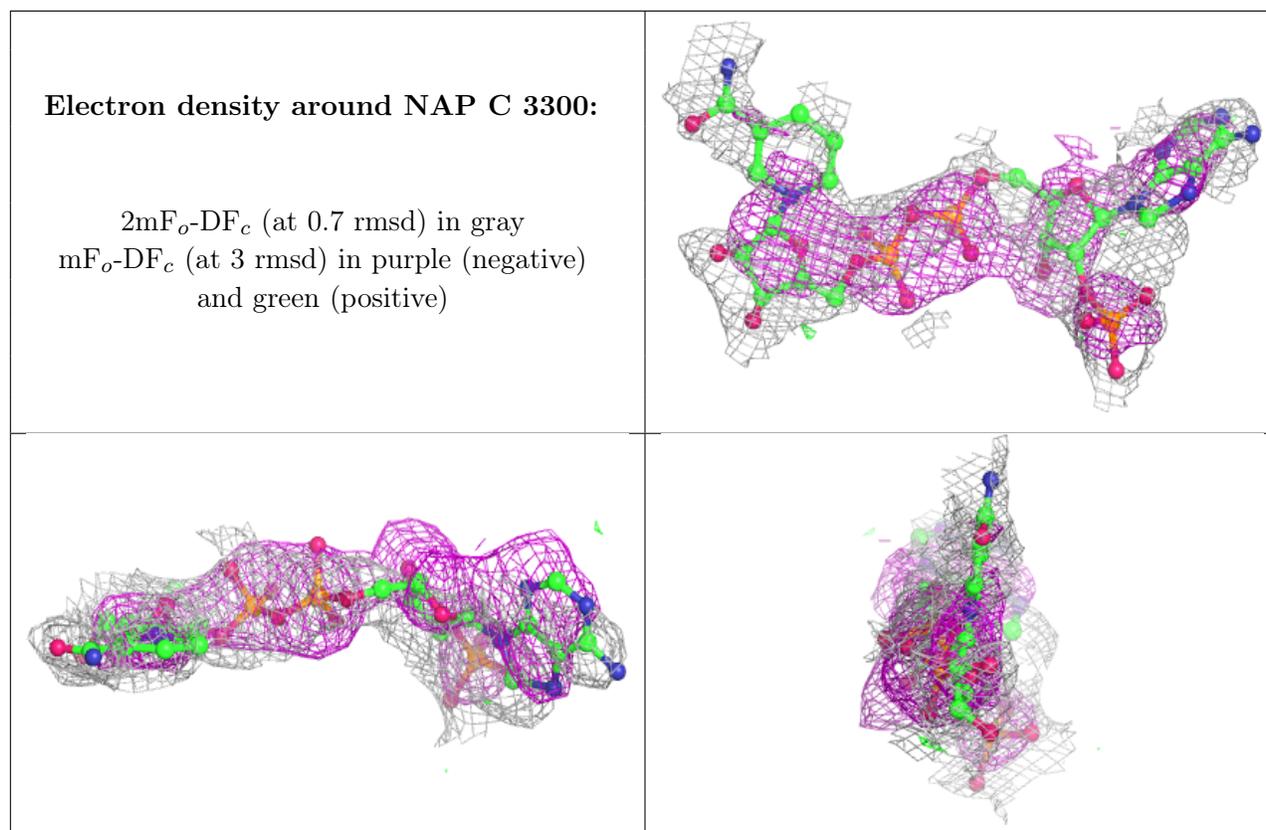
6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

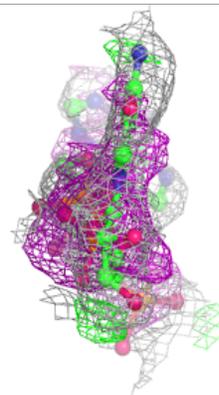
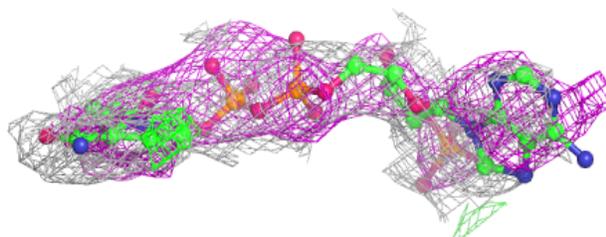
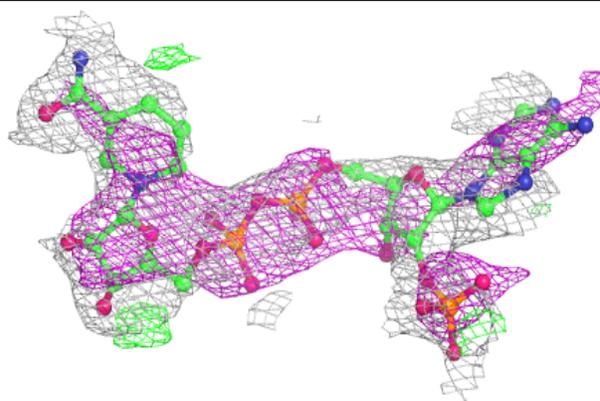
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GLU	B	2301	10/10	0.36	0.52	132,200,200,200	0
3	GLU	E	5301	10/10	0.42	0.59	157,188,190,191	0
3	GLU	C	3301	10/10	0.45	0.67	144,195,198,199	0
3	GLU	D	4301	10/10	0.55	0.51	142,156,157,158	0
3	GLU	A	1301	10/10	0.64	0.34	120,122,125,126	0
2	NAP	C	3300	48/48	0.68	0.49	79,94,101,106	0
2	NAP	D	4300	48/48	0.73	0.50	63,90,97,99	0
2	NAP	E	5300	48/48	0.74	0.50	71,90,95,99	0
2	NAP	B	2300	48/48	0.76	0.44	61,86,91,94	0
2	NAP	A	1300	48/48	0.80	0.40	55,76,86,88	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.

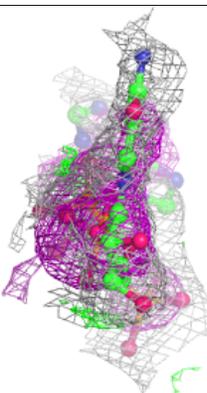
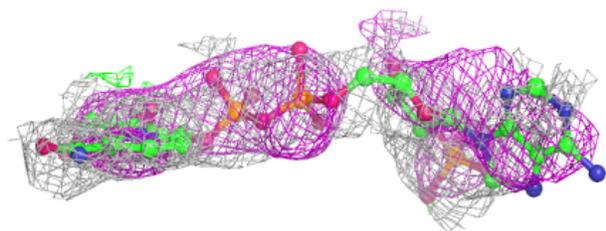
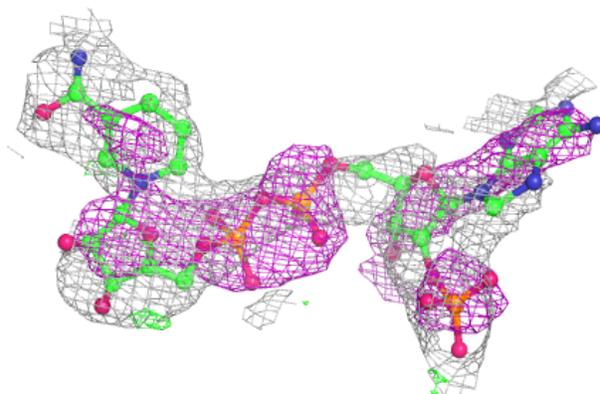


Electron density around NAP D 4300:

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

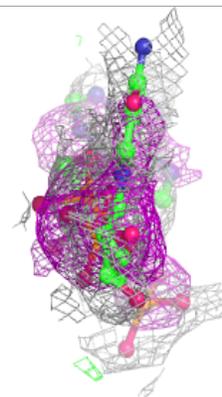
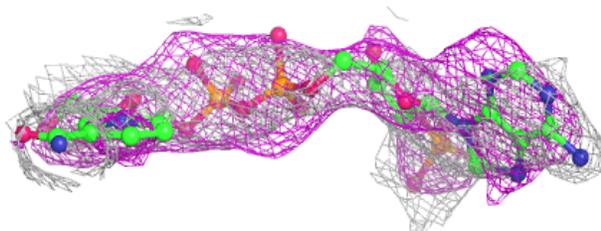
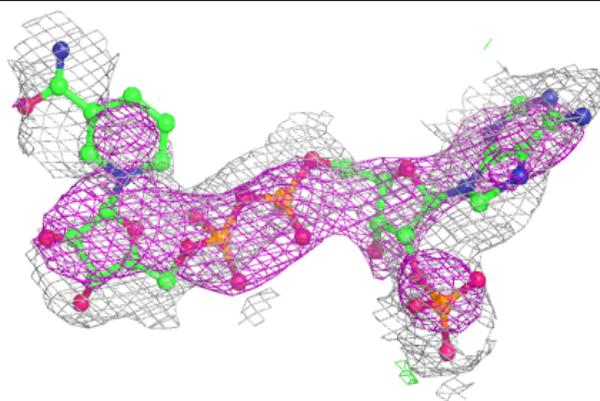
**Electron density around NAP E 5300:**

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

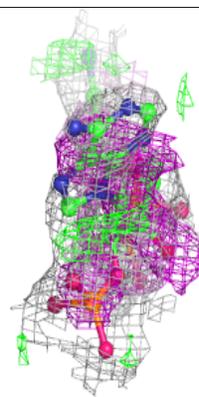
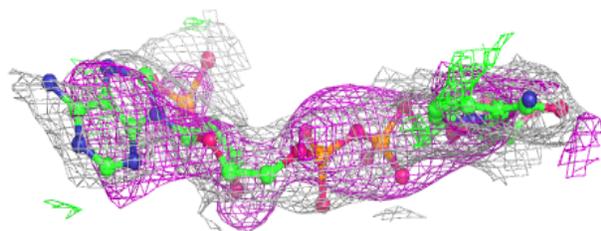
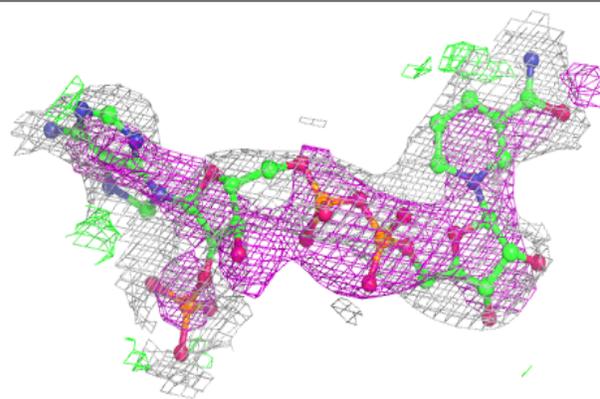


Electron density around NAP B 2300:

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

**Electron density around NAP A 1300:**

$2mF_o-DF_c$ (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.