



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 25, 2025 – 01:19 am BST

PDB ID : 9QZE / pdb\_00009qze  
Title : Exploiting ALDH1A2 and ALDH1A3 Isoform Variability for Crystallization Screening  
Authors : Garaavglia, S.; Mazzorana, M.  
Deposited on : 2025-04-22  
Resolution : 2.01 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](mailto: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>.

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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 Phenix2.0rc1
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.006 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.45.1

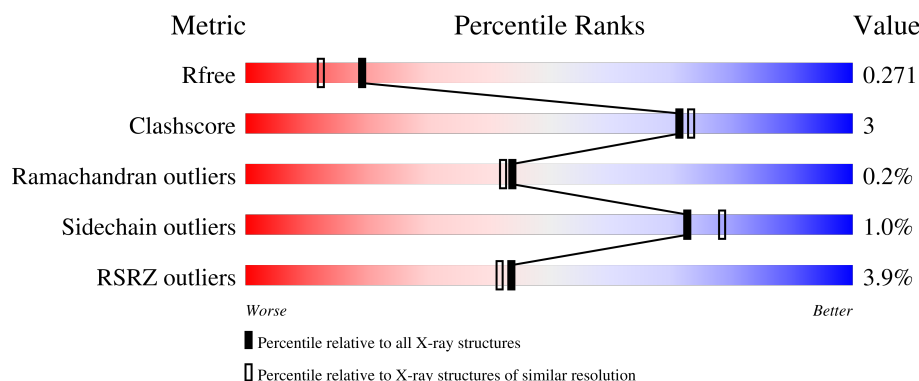
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.01 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))
$R_{free}$	164625	9409 (2.00-2.00)
Clashscore	180529	10737 (2.00-2.00)
Ramachandran outliers	177936	10628 (2.00-2.00)
Sidechain outliers	177891	10627 (2.00-2.00)
RSRZ outliers	164620	9409 (2.00-2.00)

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	512	<div> <div></div> <div>85%</div> <div>9%</div> <div>6%</div> </div>
1	B	512	<div> <div>6%</div> <div>86%</div> <div>7%</div> <div>7%</div> </div>

## 2 Entry composition [i](#)

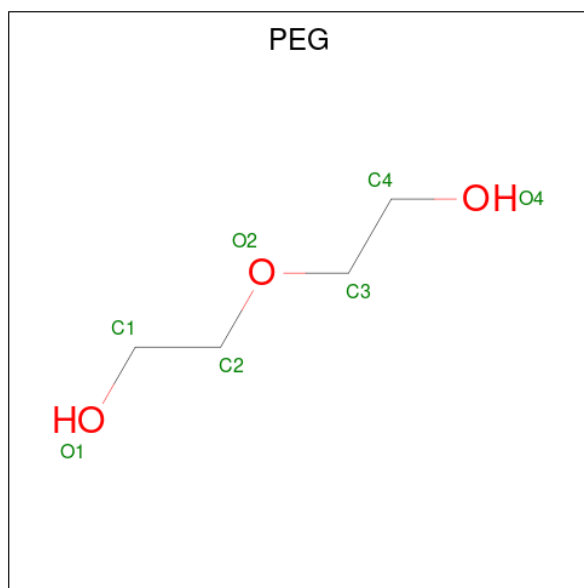
There are 4 unique types of molecules in this entry. The entry contains 7678 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Aldehyde dehydrogenase family 1 member A3.

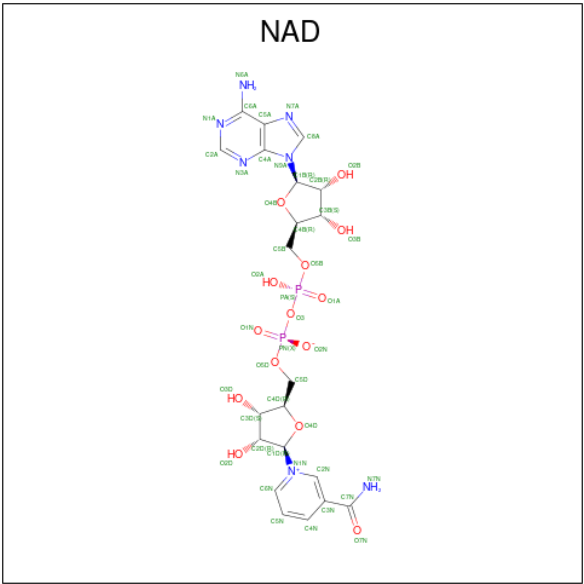
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	481	Total	C	N	O	S	0	0	0
			3707	2363	631	693	20			
1	B	478	Total	C	N	O	S	0	1	0
			3691	2353	628	689	21			

- Molecule 2 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			7	4	3		
2	B	1	Total	C	O	0	0
			7	4	3		
2	B	1	Total	C	O	0	0
			7	4	3		
2	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula:  $C_{21}H_{27}N_7O_{14}P_2$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	B	1	Total	C	N	O	P	0	0
			35	15	5	13	2		

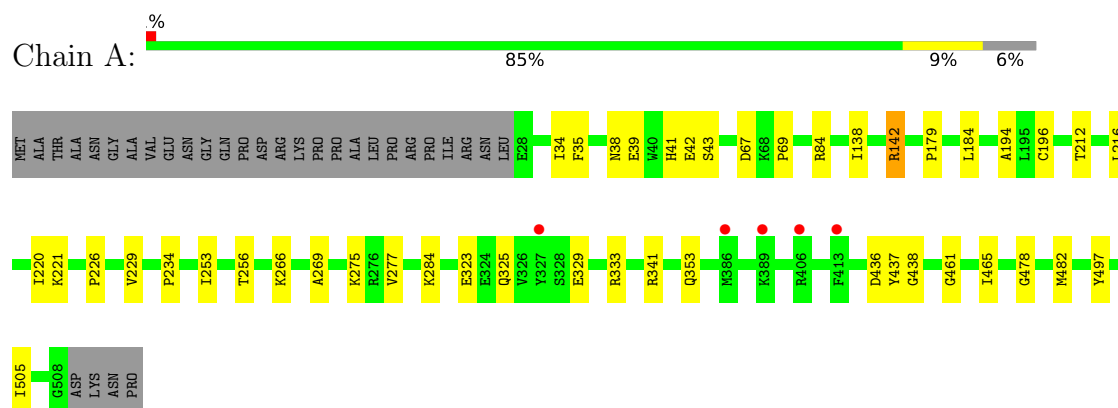
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	82	Total	O	0	0
			82	82		
4	B	91	Total	O	0	0
			91	91		

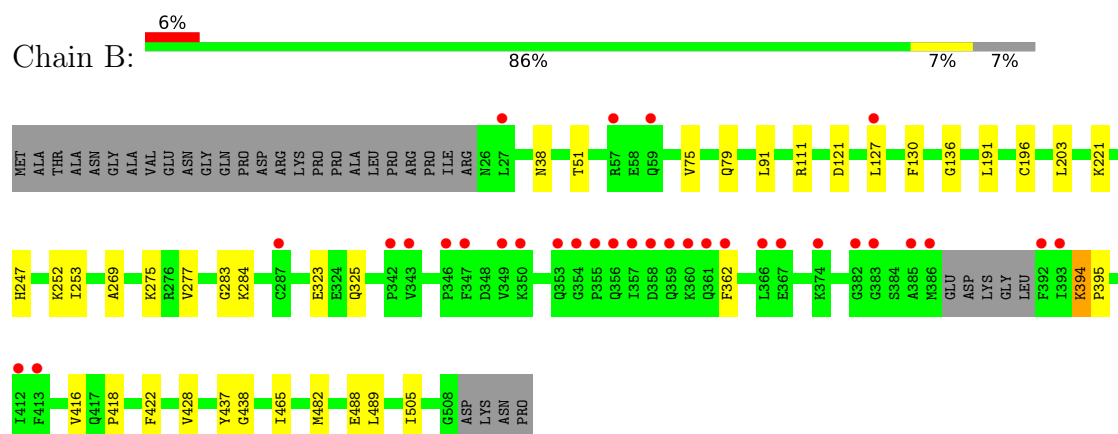
### 3 Residue-property plots [i](#)

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: Aldehyde dehydrogenase family 1 member A3



- Molecule 1: Aldehyde dehydrogenase family 1 member A3



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 <sub>1</sub> 2 <sub>1</sub> 2 <sub>1</sub>	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	81.03Å 89.20Å 158.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	79.31 – 2.01 79.31 – 2.01	Depositor EDS
% Data completeness (in resolution range)	99.8 (79.31-2.01) 99.9 (79.31-2.01)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.06 (at 2.02Å)	Xtriage
Refinement program	PHENIX 1.21.2-5419	Depositor
R, $R_{free}$	0.216 , 0.268 0.218 , 0.271	Depositor DCC
$R_{free}$ test set	3859 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.3	Xtriage
Anisotropy	0.437	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 45.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7678	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.84% 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: NAD, 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	0.39	0/3783	0.55	0/5118
1	B	0.37	0/3766	0.56	0/5095
All	All	0.38	0/7549	0.56	0/10213

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	142	ARG	Sidechain

### 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	3707	0	3728	26	0
1	B	3691	0	3709	24	0
2	A	7	0	10	0	0
2	B	21	0	30	1	0
3	A	44	0	26	0	0
3	B	35	0	19	0	0
4	A	82	0	0	0	0
4	B	91	0	0	1	0
All	All	7678	0	7522	48	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 3.

All (48) 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:329:GLU:HG3	1:A:333:ARG:HE	1.58	0.69
1:A:39:GLU:HB2	1:A:41:HIS:HE1	1.61	0.65
1:B:269:ALA:HB1	1:B:275:LYS:HG3	1.80	0.64
1:A:39:GLU:HB2	1:A:41:HIS:CE1	2.37	0.60
1:A:266:LYS:HE2	4:B:702:HOH:O	2.02	0.59
1:A:341:ARG:NH1	1:A:353:GLN:HB2	2.21	0.54
1:B:323:GLU:HG2	1:B:325:GLN:NE2	2.22	0.54
1:A:341:ARG:HH11	1:A:353:GLN:HB2	1.72	0.54
1:B:284:LYS:HE3	1:B:437:TYR:CD1	2.43	0.53
1:B:136:GLY:HA2	2:B:603:PEG:H32	1.91	0.52
1:A:465:ILE:HD12	1:B:505:ILE:HG12	1.92	0.51
1:A:323:GLU:HG2	1:A:325:GLN:OE1	2.11	0.51
1:A:329:GLU:HG2	1:A:333:ARG:HH21	1.77	0.49
1:B:284:LYS:HE3	1:B:437:TYR:CE1	2.48	0.49
1:A:269:ALA:HB1	1:A:275:LYS:HG3	1.93	0.48
1:A:253:ILE:CG2	1:A:277:VAL:HG22	2.44	0.48
1:A:184:LEU:HD13	1:A:212:THR:HB	1.95	0.47
1:B:111:ARG:HG2	1:B:130:PHE:CE2	2.50	0.47
1:B:416:VAL:O	1:B:418:PRO:HD3	2.14	0.47
1:B:325:GLN:NE2	1:B:325:GLN:H	2.13	0.47
1:A:138:ILE:O	1:A:142:ARG:HG3	2.14	0.47
1:A:179:PRO:HD3	1:A:256:THR:HB	1.97	0.46
1:B:75:VAL:HG11	1:B:247:HIS:CE1	2.51	0.46
1:A:284:LYS:HE3	1:A:437:TYR:CE1	2.51	0.45
1:A:226:PRO:O	1:A:229:VAL:HG23	2.16	0.45
1:A:436:ASP:HB3	1:A:482:MET:HE3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:461:GLY:HA3	1:A:478:GLY:O	2.17	0.44
1:B:38:ASN:HB3	1:B:221:LYS:HD2	1.99	0.44
1:A:34:ILE:HG12	1:A:234:PRO:HD2	1.99	0.43
1:B:253:ILE:HG23	1:B:277:VAL:HG13	2.00	0.43
1:B:91:LEU:HD23	1:B:91:LEU:HA	1.88	0.43
1:A:216:LEU:O	1:A:220:ILE:HG13	2.18	0.43
1:A:67:ASP:HB3	1:A:69:PRO:HD2	2.01	0.42
1:B:488:GLU:O	1:B:489:LEU:HB2	2.19	0.42
1:A:505:ILE:HG12	1:B:465:ILE:HD12	2.02	0.42
1:B:127:LEU:HD23	1:B:127:LEU:HA	1.88	0.42
1:B:253:ILE:HG22	1:B:277:VAL:HG22	2.02	0.42
1:B:283:GLY:HA2	1:B:437:TYR:HB3	2.01	0.42
1:A:38:ASN:HB3	1:A:221:LYS:HD2	2.02	0.42
1:A:194:ALA:HB2	1:A:497:TYR:CE1	2.55	0.42
1:B:191:LEU:C	1:B:191:LEU:HD23	2.45	0.41
1:A:35:PHE:CZ	1:A:38:ASN:HA	2.55	0.41
1:B:394:LYS:HD3	1:B:395:PRO:HD2	2.02	0.41
1:A:42:GLU:HG2	1:A:43:SER:O	2.21	0.41
1:B:482:MET:HE2	1:B:482:MET:HB3	1.92	0.40
1:B:362:PHE:C	1:B:362:PHE:CD2	2.99	0.40
1:B:75:VAL:O	1:B:79:GLN:HG3	2.22	0.40
1:B:422:PHE:CD2	1:B:428:VAL:HB	2.57	0.40

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	479/512 (94%)	459 (96%)	19 (4%)	1 (0%)	44	42
1	B	475/512 (93%)	460 (97%)	14 (3%)	1 (0%)	44	42
All	All	954/1024 (93%)	919 (96%)	33 (4%)	2 (0%)	44	42

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	438	GLY
1	A	438	GLY

### 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	396/422 (94%)	394 (100%)	2 (0%)	86	90
1	B	395/422 (94%)	389 (98%)	6 (2%)	60	66
All	All	791/844 (94%)	783 (99%)	8 (1%)	73	78

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	ARG
1	A	196	CYS
1	B	51	THR
1	B	121	ASP
1	B	196	CYS
1	B	203	LEU
1	B	252	LYS
1	B	394	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	128	HIS
1	A	152	GLN
1	A	356	GLN
1	A	485	ASN
1	B	312	GLN
1	B	325	GLN
1	B	353	GLN
1	B	485	ASN

### 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 ⓘ

6 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$
3	NAD	A	602	-	42,48,48	0.89	2 (4%)	50,73,73	0.86	1 (2%)
2	PEG	A	601	-	6,6,6	0.26	0	5,5,5	0.32	0
2	PEG	B	602	-	6,6,6	0.23	0	5,5,5	0.52	0
3	NAD	B	604	-	33,38,48	1.03	2 (6%)	37,58,73	1.03	2 (5%)
2	PEG	B	603	-	6,6,6	0.37	0	5,5,5	0.67	0
2	PEG	B	601	-	6,6,6	0.23	0	5,5,5	0.26	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	NAD	A	602	-	-	14/26/62/62	0/5/5/5
2	PEG	A	601	-	-	2/4/4/4	-
2	PEG	B	602	-	-	3/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAD	B	604	-	-	8/18/51/62	0/4/4/5
2	PEG	B	603	-	-	2/4/4/4	-
2	PEG	B	601	-	-	1/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	604	NAD	O4D-C1D	-2.66	1.38	1.43
3	A	602	NAD	C8A-N7A	-2.46	1.30	1.34
3	A	602	NAD	O2B-C2B	-2.19	1.37	1.43
3	B	604	NAD	C8A-N7A	-2.09	1.31	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	604	NAD	C4A-C5A-N7A	2.47	111.97	109.40
3	B	604	NAD	O2A-PA-O1A	2.46	124.38	112.24
3	A	602	NAD	O2A-PA-O1A	2.33	123.75	112.24

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	602	NAD	C5D-O5D-PN-O3
3	A	602	NAD	C5D-O5D-PN-O1N
3	A	602	NAD	O4D-C1D-N1N-C2N
3	A	602	NAD	O4D-C1D-N1N-C6N
3	A	602	NAD	C2D-C1D-N1N-C6N
3	A	602	NAD	C2N-C3N-C7N-O7N
3	A	602	NAD	C2N-C3N-C7N-N7N
3	B	604	NAD	C5D-O5D-PN-O3
3	B	604	NAD	C5D-O5D-PN-O1N
3	B	604	NAD	C5D-O5D-PN-O2N
3	B	604	NAD	O4D-C4D-C5D-O5D
3	A	602	NAD	C4N-C3N-C7N-O7N
3	A	602	NAD	C4N-C3N-C7N-N7N
2	B	603	PEG	O2-C3-C4-O4
3	B	604	NAD	C3D-C4D-C5D-O5D
2	B	601	PEG	O1-C1-C2-O2
2	A	601	PEG	O1-C1-C2-O2
2	A	601	PEG	O2-C3-C4-O4

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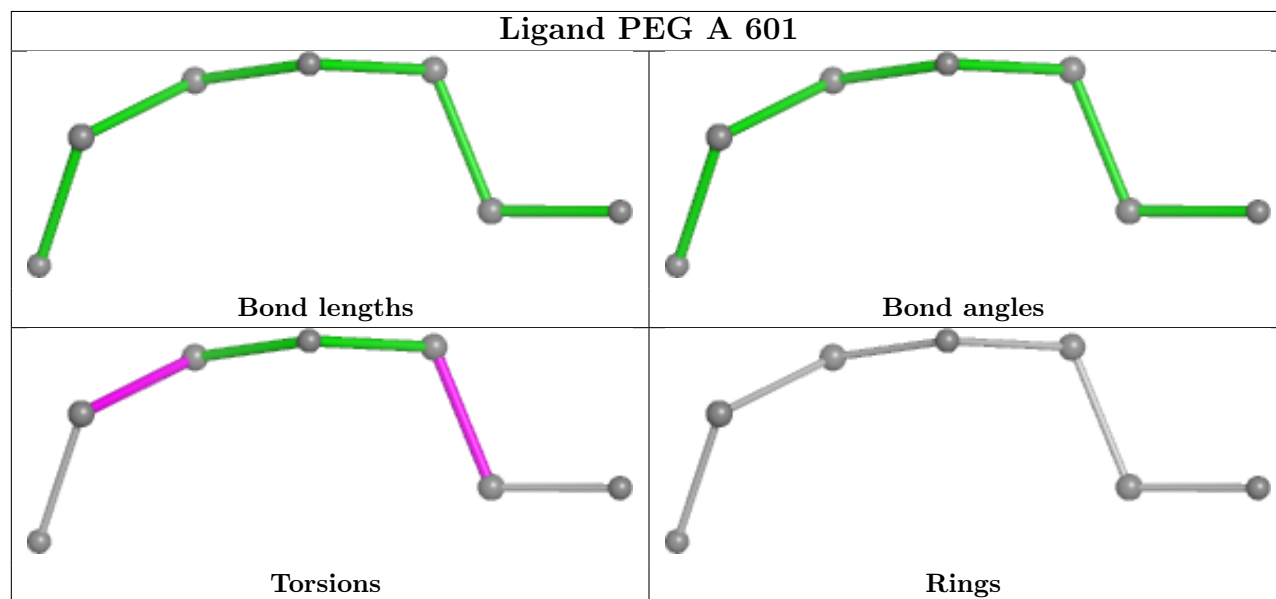
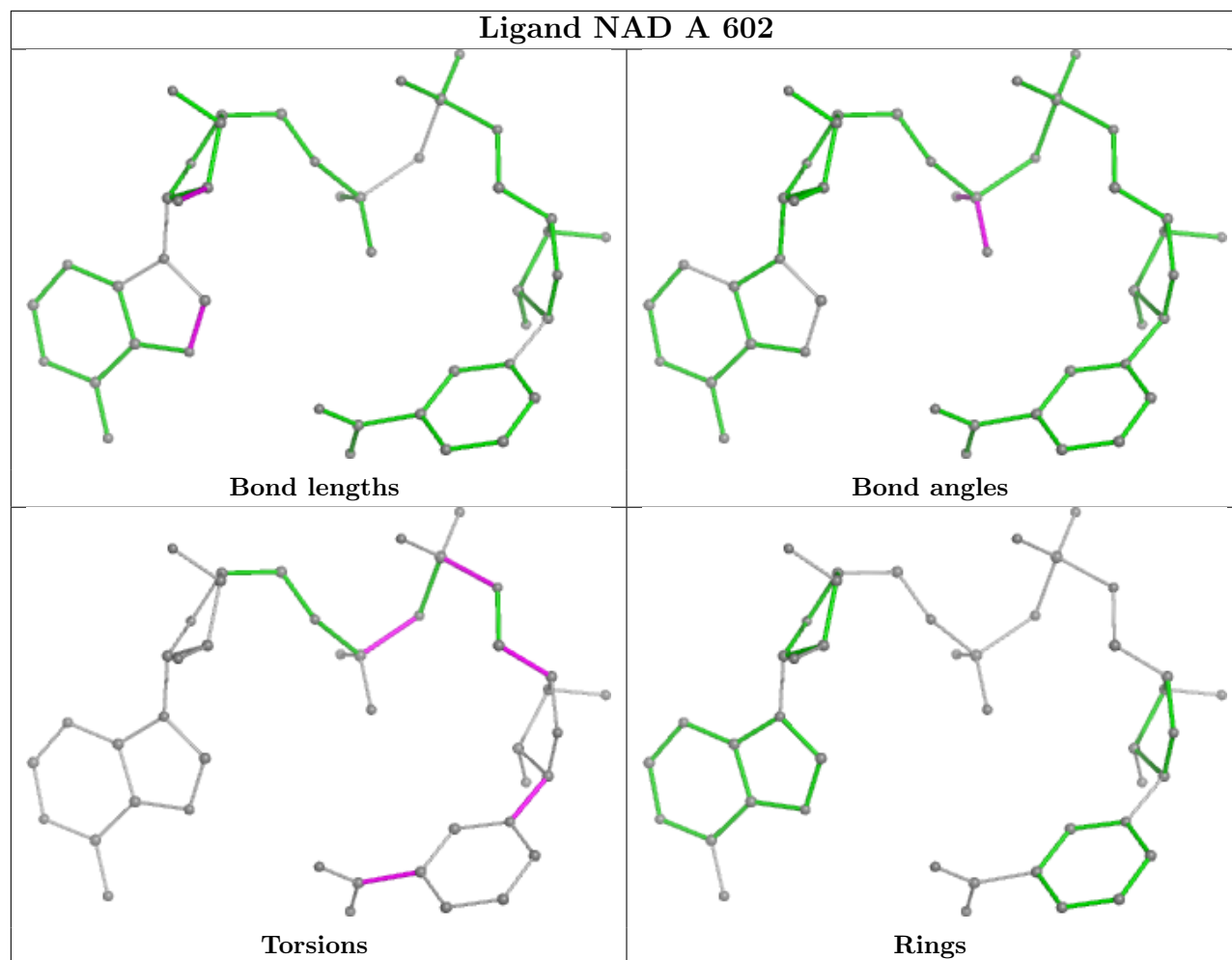
Mol	Chain	Res	Type	Atoms
2	B	602	PEG	O1-C1-C2-O2
2	B	602	PEG	O2-C3-C4-O4
3	B	604	NAD	PN-O3-PA-O1A
3	A	602	NAD	PN-O3-PA-O5B
3	A	602	NAD	C5D-O5D-PN-O2N
3	B	604	NAD	PA-O3-PN-O1N
2	B	603	PEG	C4-C3-O2-C2
3	B	604	NAD	PN-O3-PA-O5B
3	A	602	NAD	C2D-C1D-N1N-C2N
3	A	602	NAD	PN-O3-PA-O1A
3	A	602	NAD	O4D-C4D-C5D-O5D
2	B	602	PEG	C4-C3-O2-C2

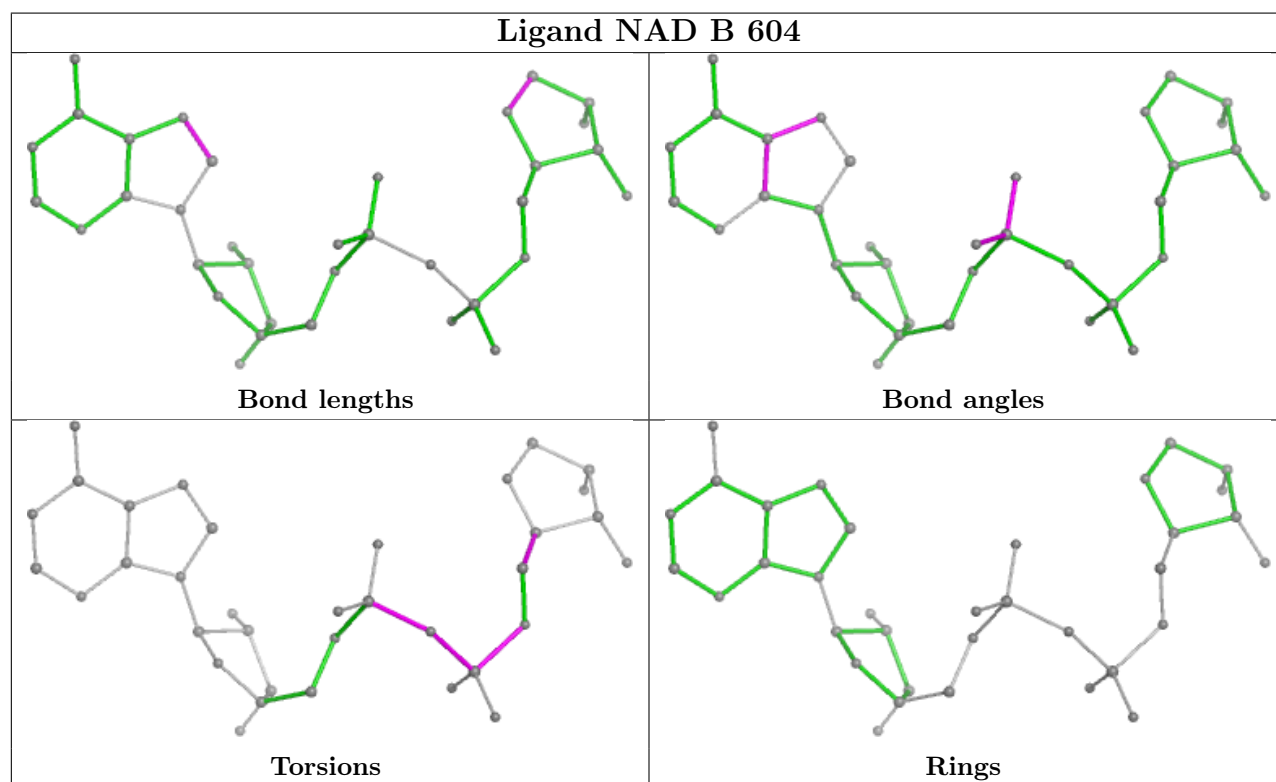
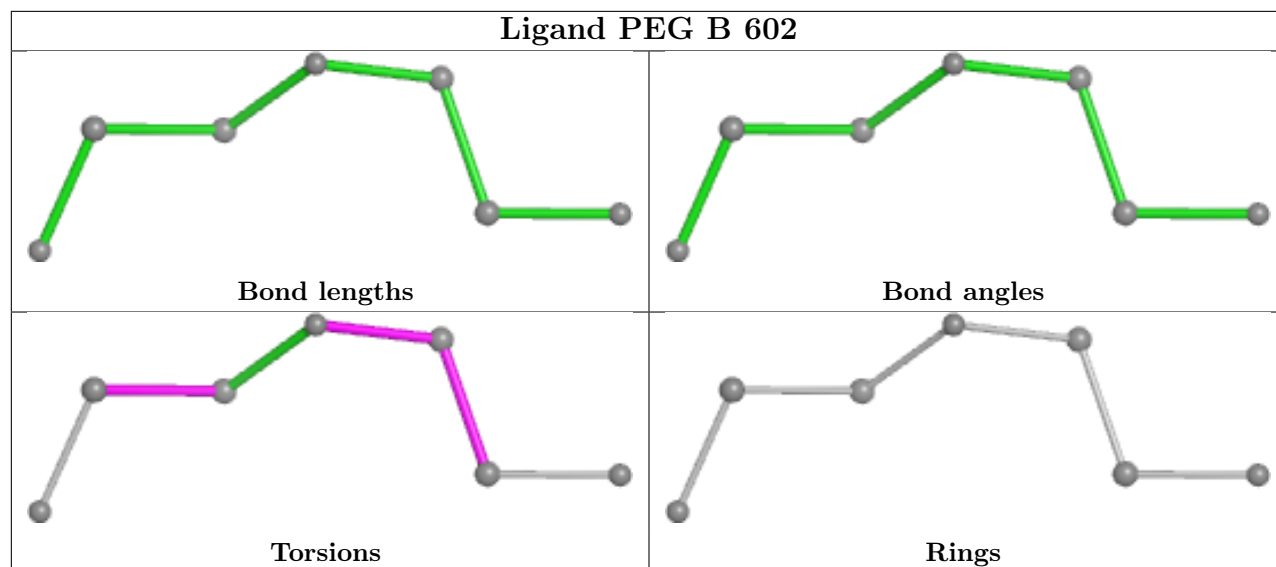
There are no ring outliers.

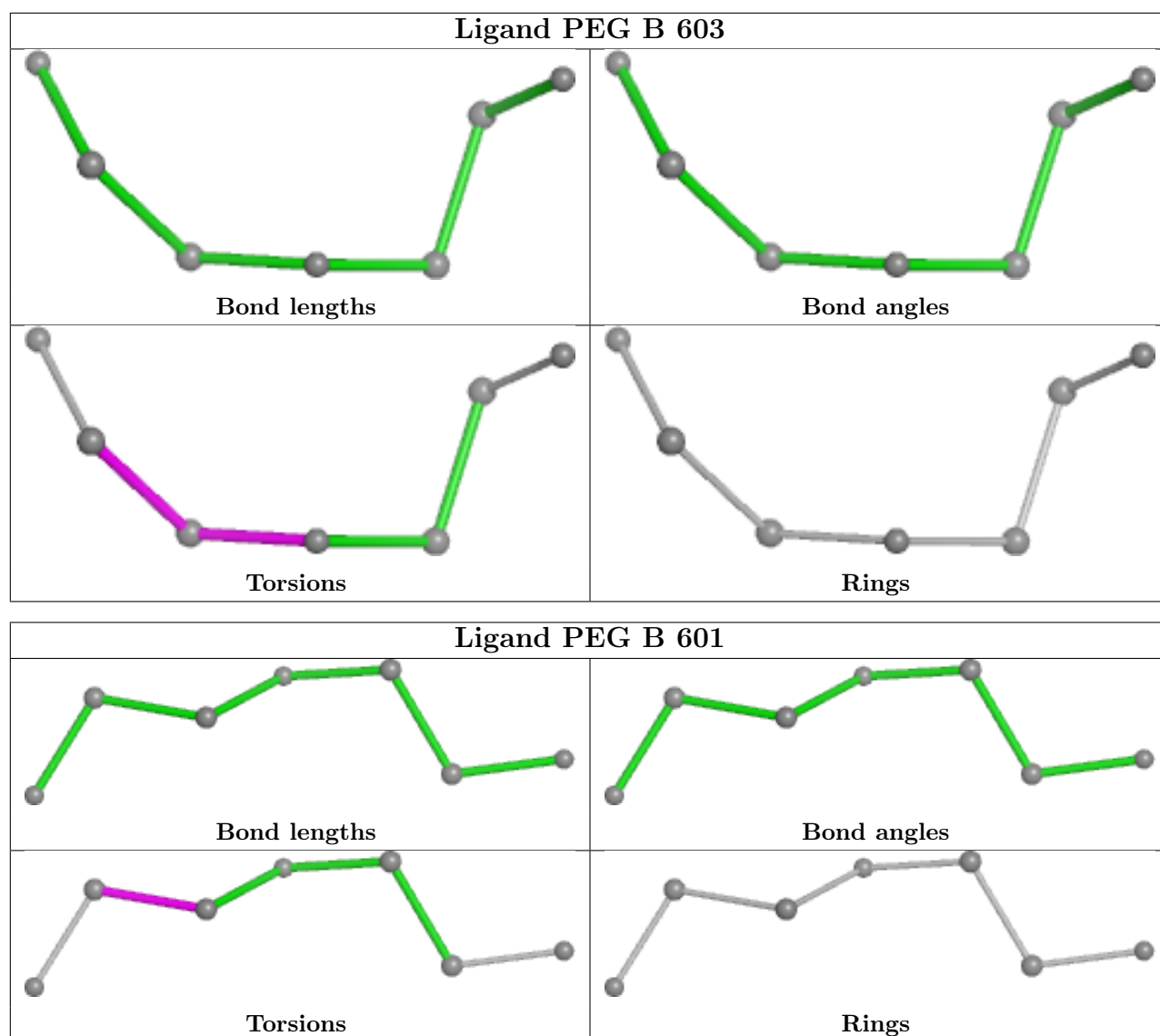
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	603	PEG	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, 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	481/512 (93%)	0.18	5 (1%) 79 78	25, 40, 64, 114	0
1	B	478/512 (93%)	0.43	32 (6%) 25 23	24, 40, 96, 153	1 (0%)
All	All	959/1024 (93%)	0.30	37 (3%) 44 42	24, 40, 80, 153	1 (0%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	357	ILE	5.9
1	B	413	PHE	5.8
1	B	392	PHE	5.3
1	B	347	PHE	5.1
1	B	358	ASP	4.4
1	B	361	GLN	4.3
1	B	349	VAL	3.9
1	B	393	ILE	3.6
1	A	413	PHE	3.6
1	B	356	GLN	3.4
1	B	343	VAL	3.4
1	B	382	GLY	3.2
1	B	57	ARG	3.2
1	B	386	MET	3.1
1	B	342	PRO	3.0
1	B	354	GLY	3.0
1	B	355	PRO	3.0
1	B	362	PHE	3.0
1	B	360	LYS	3.0
1	B	359	GLN	2.8
1	A	389	LYS	2.8
1	B	346	PRO	2.8
1	B	27	LEU	2.7
1	A	386	MET	2.7

*Continued on next page...*

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Mol	Chain	Res	Type	RSRZ
1	B	353	GLN	2.5
1	B	350	LYS	2.5
1	B	385	ALA	2.5
1	B	287[A]	CYS	2.4
1	B	383	GLY	2.4
1	A	327	TYR	2.3
1	B	127	LEU	2.3
1	B	367	GLU	2.2
1	B	59	GLN	2.2
1	A	406	ARG	2.2
1	B	366	LEU	2.2
1	B	374	LYS	2.1
1	B	412	ILE	2.1

## 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](#)

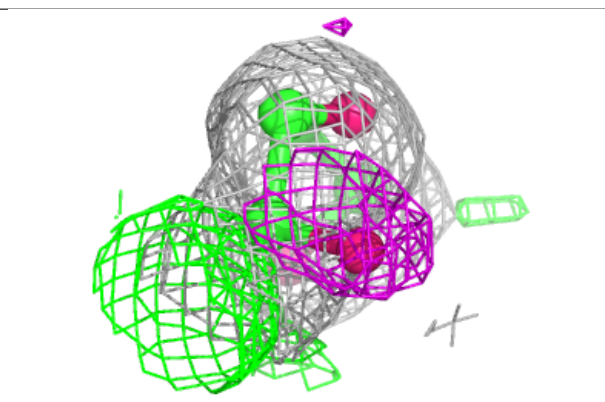
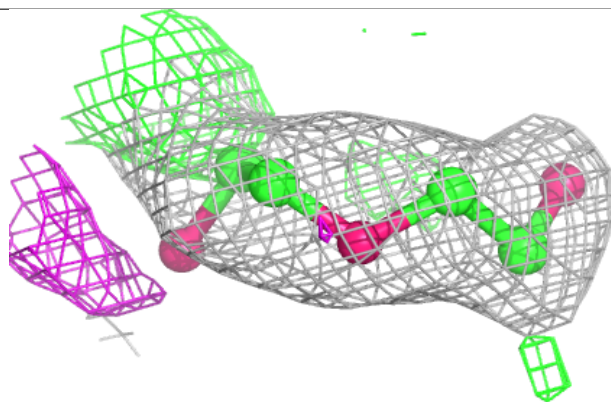
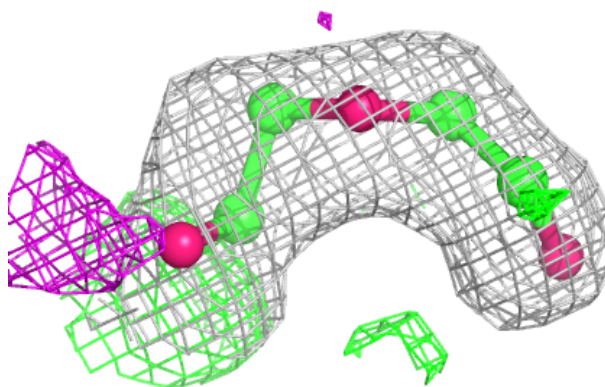
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	PEG	B	603	7/7	0.77	0.18	39,40,50,69	0
2	PEG	B	602	7/7	0.89	0.14	51,53,57,58	0
2	PEG	A	601	7/7	0.91	0.10	32,35,44,47	0
2	PEG	B	601	7/7	0.92	0.12	45,58,60,61	0
3	NAD	A	602	44/44	0.92	0.13	33,47,137,152	0
3	NAD	B	604	35/44	0.93	0.11	32,38,96,105	0

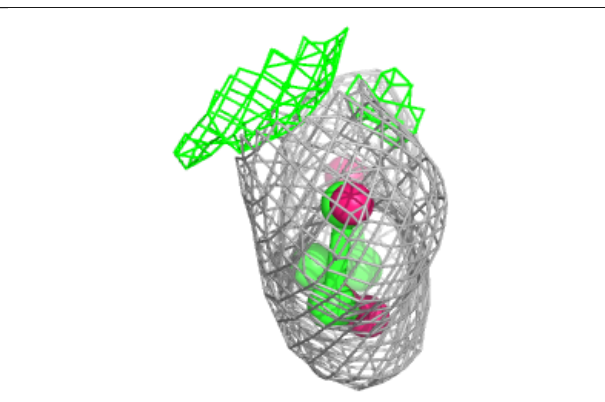
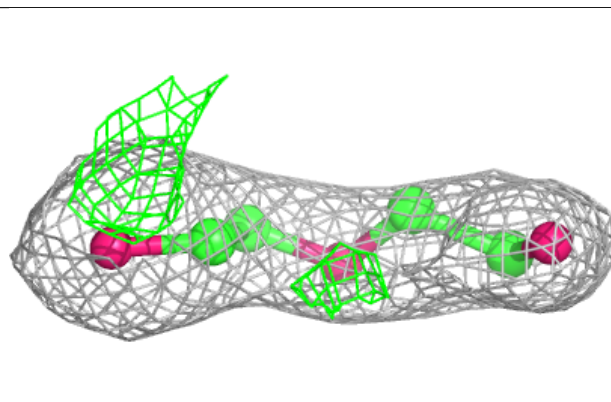
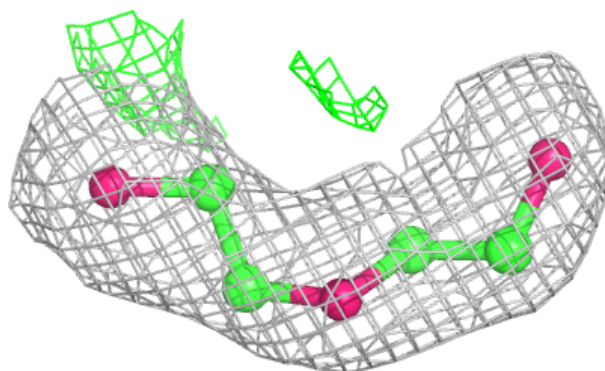
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around PEG B 603:**

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

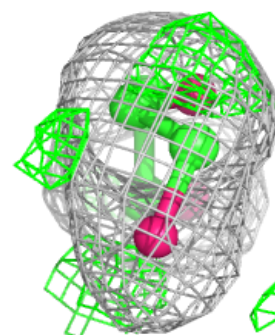
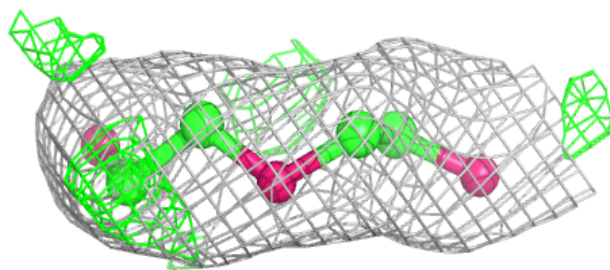
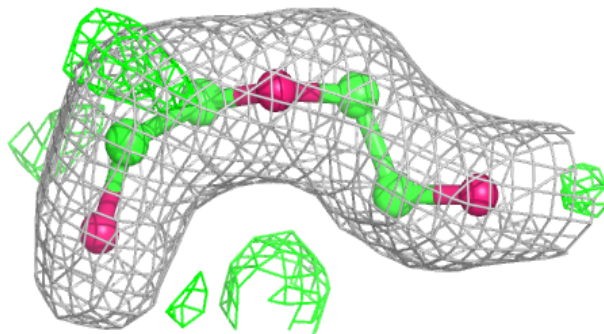
**Electron density around PEG B 602:**

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

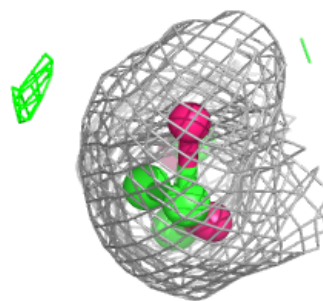
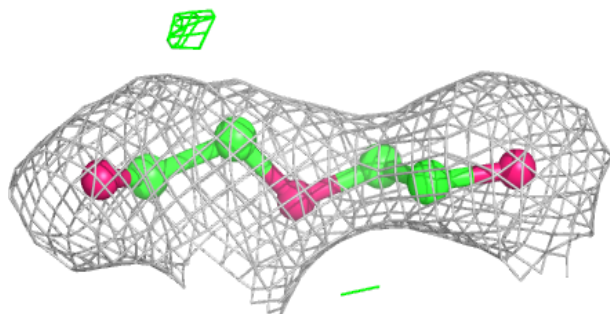
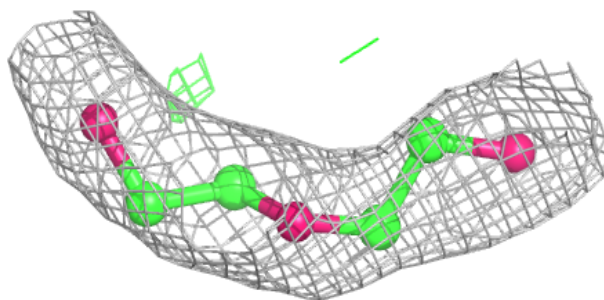


**Electron density around PEG A 601:**

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

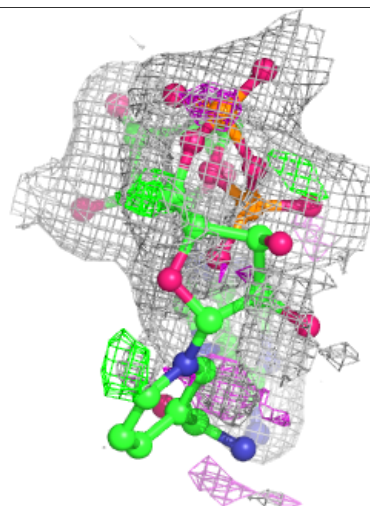
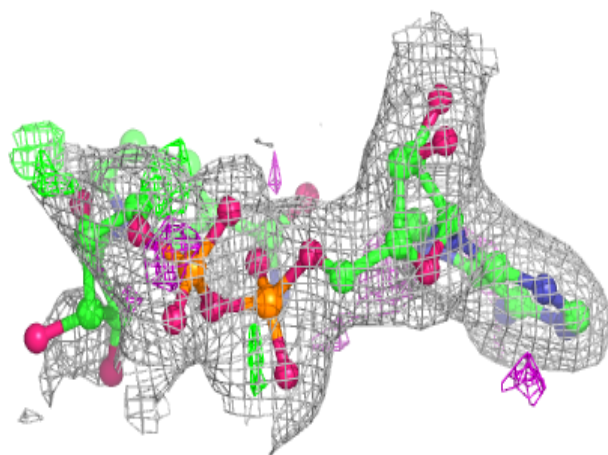
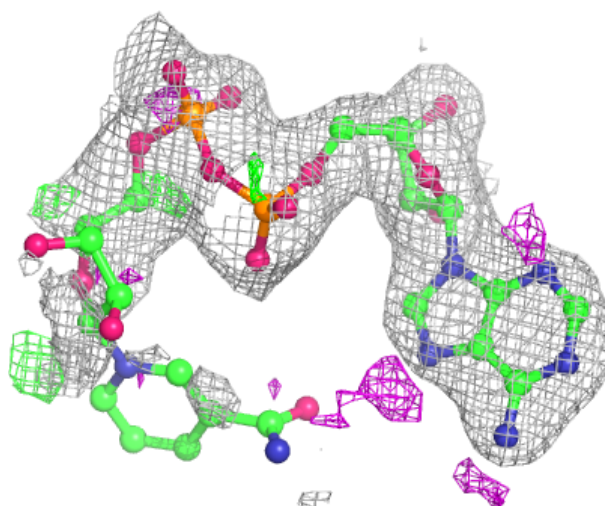
**Electron density around PEG B 601:**

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

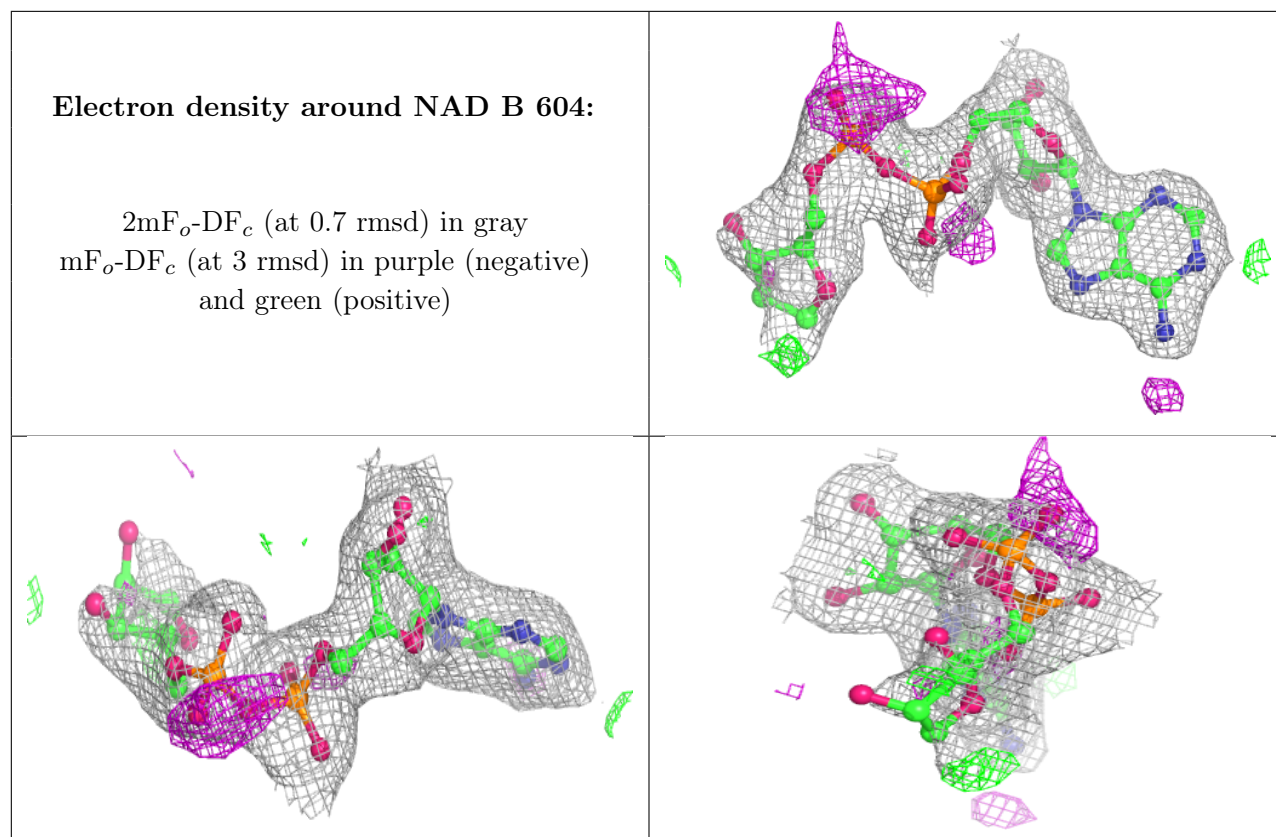


**Electron density around NAD A 602:**

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







## 6.5 Other polymers ⓘ

There are no such residues in this entry.