



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 13, 2026 – 04:14 pm BST

PDB ID : 28JK / pdb_000028jk
Title : CRYSTAL STRUCTURE OF RAT PEROXISOMAL MULTIFUNCTIONAL ENZYME TYPE-1 COMPLEXED WITH 3R-HYDROXY-4E-DECENOYL-CoA AND REDUCED NICOTINAMIDE ADENINE DINUCLEOTIDE
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Deposited on : 2026-02-05
Resolution : 2.44 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	1.8.4, CSD as541be (2020)
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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

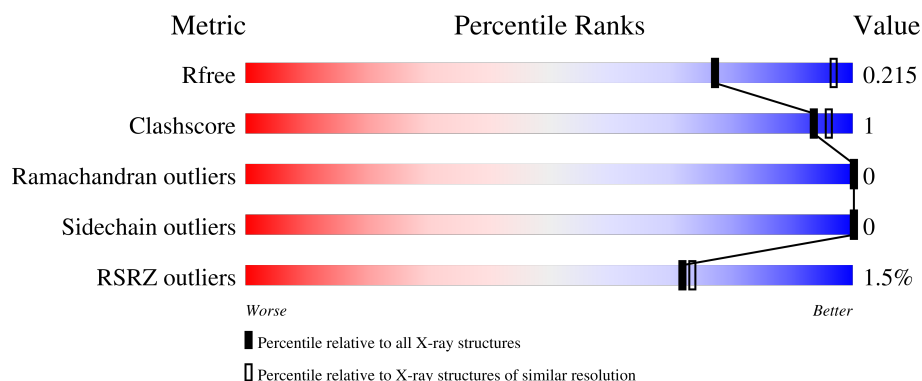
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

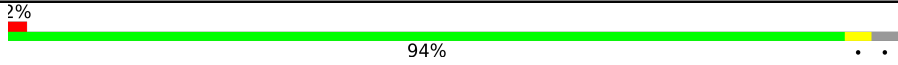
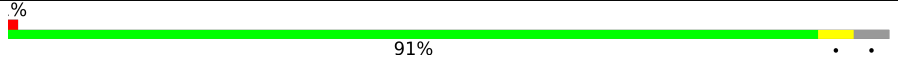
The reported resolution of this entry is 2.44 Å.

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	2340 (2.46-2.42)
Clashscore	190562	2400 (2.46-2.42)
Ramachandran outliers	187476	2379 (2.46-2.42)
Sidechain outliers	187428	2379 (2.46-2.42)
RSRZ outliers	180081	2340 (2.46-2.42)

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	742	
1	B	742	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11384 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 Peroxisomal bifunctional enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	718	Total	C	N	O	S	0	1	0
			5522	3531	967	1001	23			
1	B	710	Total	C	N	O	S	0	0	0
			5460	3495	953	989	23			

There are 42 discrepancies between the modelled and reference sequences:

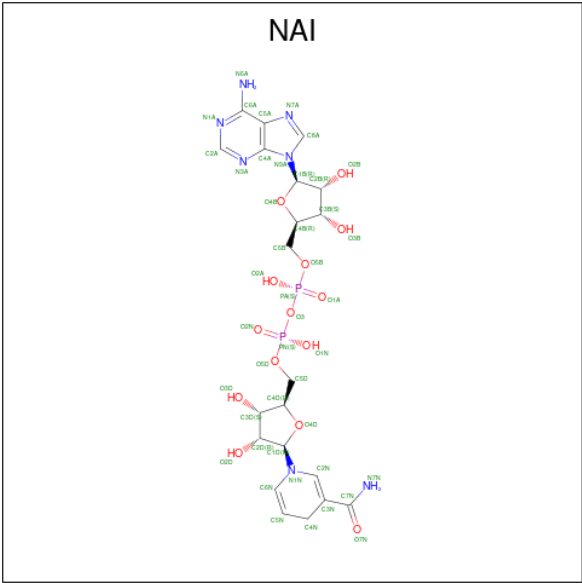
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP P07896
A	-18	GLY	-	expression tag	UNP P07896
A	-17	SER	-	expression tag	UNP P07896
A	-16	SER	-	expression tag	UNP P07896
A	-15	HIS	-	expression tag	UNP P07896
A	-14	HIS	-	expression tag	UNP P07896
A	-13	HIS	-	expression tag	UNP P07896
A	-12	HIS	-	expression tag	UNP P07896
A	-11	HIS	-	expression tag	UNP P07896
A	-10	HIS	-	expression tag	UNP P07896
A	-9	SER	-	expression tag	UNP P07896
A	-8	SER	-	expression tag	UNP P07896
A	-7	GLY	-	expression tag	UNP P07896
A	-6	LEU	-	expression tag	UNP P07896
A	-5	VAL	-	expression tag	UNP P07896
A	-4	PRO	-	expression tag	UNP P07896
A	-3	ARG	-	expression tag	UNP P07896
A	-2	GLY	-	expression tag	UNP P07896
A	-1	SER	-	expression tag	UNP P07896
A	0	HIS	-	expression tag	UNP P07896
A	514	ARG	LYS	cloning artifact	UNP P07896
B	-19	MET	-	initiating methionine	UNP P07896
B	-18	GLY	-	expression tag	UNP P07896
B	-17	SER	-	expression tag	UNP P07896
B	-16	SER	-	expression tag	UNP P07896

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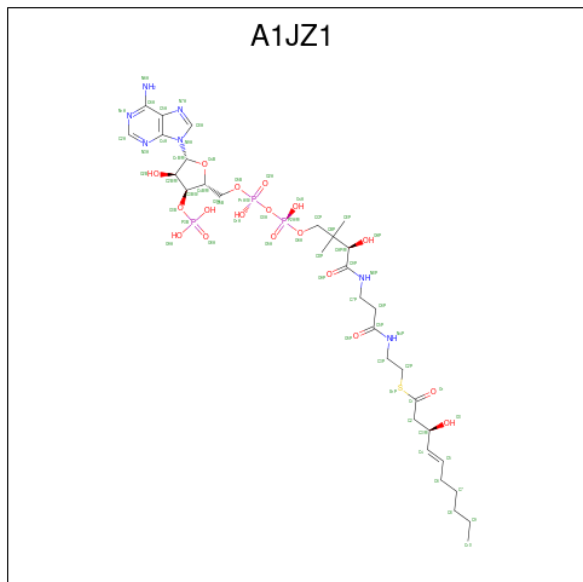
Chain	Residue	Modelled	Actual	Comment	Reference
B	-15	HIS	-	expression tag	UNP P07896
B	-14	HIS	-	expression tag	UNP P07896
B	-13	HIS	-	expression tag	UNP P07896
B	-12	HIS	-	expression tag	UNP P07896
B	-11	HIS	-	expression tag	UNP P07896
B	-10	HIS	-	expression tag	UNP P07896
B	-9	SER	-	expression tag	UNP P07896
B	-8	SER	-	expression tag	UNP P07896
B	-7	GLY	-	expression tag	UNP P07896
B	-6	LEU	-	expression tag	UNP P07896
B	-5	VAL	-	expression tag	UNP P07896
B	-4	PRO	-	expression tag	UNP P07896
B	-3	ARG	-	expression tag	UNP P07896
B	-2	GLY	-	expression tag	UNP P07896
B	-1	SER	-	expression tag	UNP P07896
B	0	HIS	-	expression tag	UNP P07896
B	514	ARG	LYS	cloning artifact	UNP P07896

- Molecule 2 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (CCD ID: NAI) (formula: C₂₁H₂₉N₇O₁₄P₂).



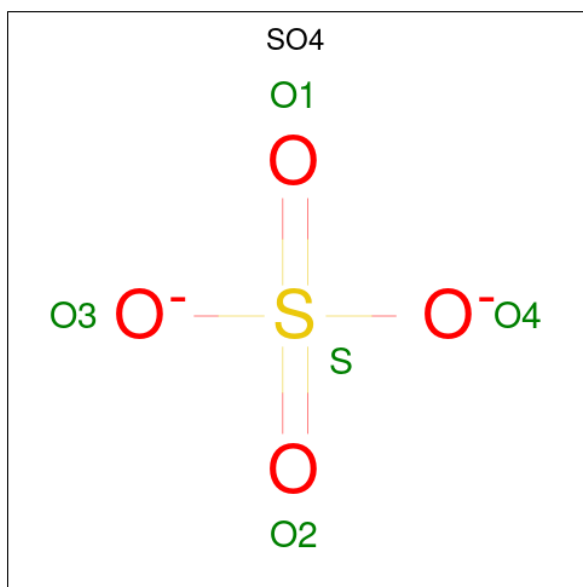
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	B	1	Total 44	C 21	N 7	O 14	P 2	0	0

- Molecule 3 is (S)-3-HYDROXY-(E)-4-DECENOYL-COA (CCD ID: A1JZ1) (formula: $C_{31}H_{52}N_7O_{18}P_3S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	0	0
			60	31	7	18	3	1		
3	B	1	Total	C	N	O	P	S	0	0
			60	31	7	18	3	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O_4S).

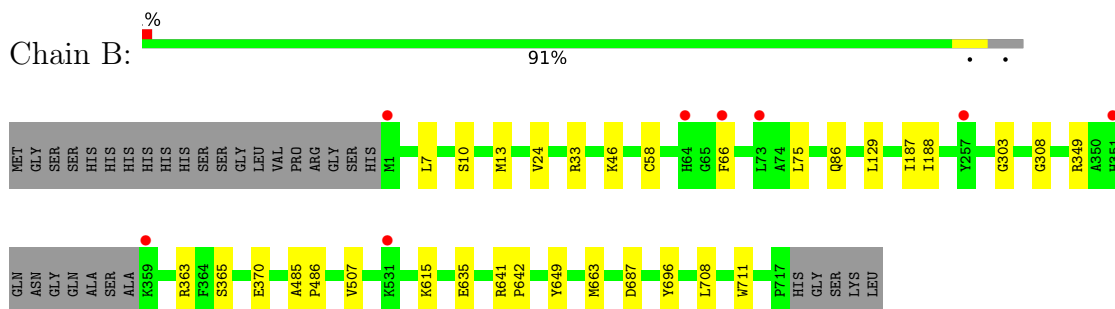


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	144	Total	O	0	0
			144	144		
5	B	35	Total	O	0	0
			35	35		

- Molecule 1: Peroxisomal bifunctional enzyme



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.28Å 126.87Å 225.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.27 – 2.44 49.27 – 2.44	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.27-2.44) 99.9 (49.27-2.44)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.174 , 0.207 0.183 , 0.215	Depositor DCC
R_{free} test set	4241 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	52.9	Xtriage
Anisotropy	0.246	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11384	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, A1JZ1, NAI

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	0.57	0/5652	0.99	1/7655 (0.0%)
1	B	0.54	0/5585	0.97	3/7566 (0.0%)
All	All	0.56	0/11237	0.98	4/15221 (0.0%)

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	2
1	B	0	2
All	All	0	4

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	200	LYS	CB-CA-C	6.76	119.11	109.38
1	B	66	PHE	CB-CA-C	-5.23	101.88	110.72
1	B	649	TYR	CA-CB-CG	-5.19	104.55	113.90
1	B	687	ASP	CA-CB-CG	5.15	117.75	112.60

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	33	ARG	Sidechain
1	A	698	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	B	33	ARG	Sidechain
1	B	349	ARG	Sidechain

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	5522	0	5634	14	0
1	B	5460	0	5576	17	0
2	A	44	0	27	0	0
2	B	44	0	27	0	0
3	A	60	0	0	0	0
3	B	60	0	0	0	0
4	A	10	0	0	0	0
4	B	5	0	0	0	0
5	A	144	0	0	1	0
5	B	35	0	0	1	0
All	All	11384	0	11264	31	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 (31) 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:46:LYS:HB2	1:B:188:ILE:HD11	1.74	0.69
1:A:515:MET:HE3	1:A:519:ARG:HB3	1.82	0.60
1:A:129:LEU:C	1:A:129:LEU:HD12	2.29	0.58
1:A:9:HIS:HE1	5:A:1028:HOH:O	1.87	0.57
1:B:129:LEU:HD12	1:B:129:LEU:C	2.30	0.57
1:B:365:SER:OG	1:B:370:GLU:HG3	2.10	0.51
1:A:515:MET:HE1	1:A:523:LEU:HD12	1.92	0.51
1:B:663:MET:HA	1:B:663:MET:HE2	1.93	0.51
1:A:7:LEU:HB2	1:A:11:LEU:HB2	1.94	0.50
1:A:663:MET:HA	1:A:663:MET:HE2	1.94	0.48
1:B:24:VAL:HG23	1:B:58:CYS:SG	2.53	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:303:GLY:O	1:B:308:GLY:HA3	2.14	0.48
1:B:635:GLU:HB3	1:B:696:TYR:HB2	1.96	0.47
1:B:10:SER:HB3	1:B:46:LYS:HG2	1.97	0.46
1:A:365:SER:OG	1:A:370:GLU:HG3	2.16	0.46
1:A:187:ILE:C	1:A:187:ILE:HD12	2.40	0.45
1:B:187:ILE:C	1:B:187:ILE:HD12	2.41	0.45
1:A:303:GLY:O	1:A:308:GLY:HA3	2.16	0.45
1:A:494:LEU:HD22	1:A:613:ILE:HG21	1.98	0.45
1:B:86:GLN:NE2	5:B:904:HOH:O	2.51	0.44
1:B:7:LEU:HD11	1:B:13:MET:HE3	2.00	0.44
1:B:24:VAL:HG11	1:B:75:LEU:HD13	2.00	0.44
1:B:485:ALA:HB3	1:B:486:PRO:HD3	2.01	0.43
1:B:507:VAL:HG22	1:B:615:LYS:HD2	2.01	0.42
1:B:363:ARG:HH12	1:B:370:GLU:HB3	1.85	0.42
1:A:7:LEU:HB3	1:A:8:PRO:HD2	2.02	0.42
1:B:641:ARG:HB2	1:B:642:PRO:HD2	2.02	0.41
1:A:484:LEU:HD21	1:A:488:TYR:HE2	1.85	0.41
1:A:7:LEU:HD11	1:A:13:MET:HE3	2.03	0.40
1:A:345:LYS:O	1:A:349:ARG:HG2	2.21	0.40
1:B:708:LEU:HA	1:B:711:TRP:CE2	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	715/742 (96%)	700 (98%)	15 (2%)	0	100	100
1	B	706/742 (95%)	690 (98%)	16 (2%)	0	100	100
All	All	1421/1484 (96%)	1390 (98%)	31 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	591/609 (97%)	591 (100%)	0	100	100
1	B	584/609 (96%)	584 (100%)	0	100	100
All	All	1175/1218 (96%)	1175 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	474	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	B	803	-	4,4,4	0.34	0	6,6,6	0.17	0
2	NAI	A	801	-	45,48,48	0.49	0	60,73,73	0.66	0
4	SO4	A	804	-	4,4,4	0.31	0	6,6,6	0.09	0
3	A1JZ1	A	802	-	56,62,62	2.02	13 (23%)	75,89,89	2.13	22 (29%)
4	SO4	A	803	-	4,4,4	0.24	0	6,6,6	0.32	0
3	A1JZ1	B	802	-	56,62,62	1.51	7 (12%)	75,89,89	1.95	17 (22%)
2	NAI	B	801	-	45,48,48	0.49	0	60,73,73	0.73	1 (1%)

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	A1JZ1	B	802	-	-	14/62/78/78	0/3/3/3
2	NAI	B	801	-	-	4/29/72/72	0/5/5/5
2	NAI	A	801	-	-	7/29/72/72	0/5/5/5
3	A1JZ1	A	802	-	-	12/62/78/78	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	802	A1JZ1	C1-S1P	9.01	1.98	1.76
3	B	802	A1JZ1	C4-C5	5.21	1.54	1.31
3	A	802	A1JZ1	C4-C5	4.96	1.52	1.31
3	A	802	A1JZ1	C2A-N1A	4.68	1.42	1.33
3	B	802	A1JZ1	C2A-N1A	3.93	1.41	1.33
3	A	802	A1JZ1	C2A-N3A	3.89	1.41	1.33
3	B	802	A1JZ1	C2A-N3A	3.34	1.40	1.33
3	A	802	A1JZ1	O3-C3	2.89	1.49	1.43
3	A	802	A1JZ1	C5A-N7A	-2.67	1.34	1.39
3	A	802	A1JZ1	C3-C4	2.65	1.55	1.50
3	B	802	A1JZ1	C3-C4	2.58	1.55	1.50
3	A	802	A1JZ1	C8A-N9A	-2.42	1.33	1.37
3	B	802	A1JZ1	C1-S1P	2.27	1.81	1.76
3	A	802	A1JZ1	O1-C1	2.24	1.24	1.21
3	A	802	A1JZ1	C2-C3	2.20	1.57	1.54
3	B	802	A1JZ1	C5A-C4A	-2.17	1.35	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	802	A1JZ1	C5A-C4A	-2.09	1.35	1.39
3	B	802	A1JZ1	C8A-N9A	-2.09	1.33	1.37
3	A	802	A1JZ1	C4A-N9A	-2.07	1.33	1.37
3	A	802	A1JZ1	O4B-C1B	2.04	1.46	1.42

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	802	A1JZ1	N3A-C2A-N1A	-7.26	117.24	128.60
3	B	802	A1JZ1	N3A-C2A-N1A	-6.89	117.83	128.60
3	A	802	A1JZ1	O4B-C1B-N9A	4.70	117.32	108.06
3	B	802	A1JZ1	N9A-C8A-N7A	-4.65	107.55	113.91
3	A	802	A1JZ1	N9A-C8A-N7A	-4.45	107.83	113.91
3	B	802	A1JZ1	O6A-CCP-CBP	-4.17	103.85	110.55
3	A	802	A1JZ1	O1-C1-S1P	-4.15	117.22	122.61
3	B	802	A1JZ1	C6P-C7P-N8P	-4.13	103.56	111.90
3	A	802	A1JZ1	C7P-N8P-C9P	4.11	129.92	122.59
3	B	802	A1JZ1	C5A-C4A-N3A	-4.03	121.49	126.75
3	A	802	A1JZ1	O6A-CCP-CBP	-3.85	104.36	110.55
3	A	802	A1JZ1	C5A-C4A-N3A	-3.83	121.75	126.75
3	B	802	A1JZ1	C7P-C6P-C5P	-3.72	106.16	112.36
3	A	802	A1JZ1	C2P-C3P-N4P	-3.67	104.70	112.42
3	A	802	A1JZ1	C5A-N7A-C8A	3.66	108.72	103.51
3	B	802	A1JZ1	C5A-N7A-C8A	3.49	108.47	103.51
3	A	802	A1JZ1	O3-C3-C2	3.47	118.79	109.73
3	B	802	A1JZ1	C2A-N3A-C4A	3.45	119.91	111.75
3	A	802	A1JZ1	C2A-N3A-C4A	3.28	119.50	111.75
3	B	802	A1JZ1	C4A-N9A-C8A	3.17	109.17	105.73
3	A	802	A1JZ1	C2-C1-S1P	2.98	117.31	113.63
3	A	802	A1JZ1	CEP-CBP-CCP	-2.94	103.43	108.23
3	B	802	A1JZ1	C5A-C6A-N6A	-2.88	117.17	123.43
3	A	802	A1JZ1	C4A-C5A-N7A	-2.83	107.17	110.62
3	A	802	A1JZ1	C6P-C7P-N8P	-2.79	106.26	111.90
3	A	802	A1JZ1	C4A-N9A-C8A	2.79	108.75	105.73
3	A	802	A1JZ1	C6A-C5A-C4A	2.78	120.93	117.18
3	B	802	A1JZ1	C2P-C3P-N4P	-2.77	106.59	112.42
3	B	802	A1JZ1	P2A-O3A-P1A	-2.67	123.65	132.83
3	B	802	A1JZ1	C4A-C5A-N7A	-2.57	107.48	110.62
3	B	802	A1JZ1	C7P-N8P-C9P	2.49	127.03	122.59
3	B	802	A1JZ1	N3A-C4A-N9A	2.49	131.18	127.08
3	A	802	A1JZ1	P2A-O3A-P1A	-2.48	124.32	132.83
3	A	802	A1JZ1	C5A-C6A-N6A	-2.46	118.07	123.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	802	A1JZ1	O1-C1-S1P	-2.37	119.53	122.61
3	A	802	A1JZ1	N3A-C4A-N9A	2.20	130.72	127.08
2	B	801	NAI	O2A-PA-O1A	2.06	122.43	112.24
3	B	802	A1JZ1	C5A-C6A-N1A	2.04	121.78	117.39
3	A	802	A1JZ1	CDP-CBP-CCP	2.02	111.52	108.23
3	A	802	A1JZ1	C6-C5-C4	-2.01	116.40	125.39

There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	802	A1JZ1	C1-C2-C3-O3
3	A	802	A1JZ1	C5B-O5B-P1A-O1A
3	A	802	A1JZ1	C5B-O5B-P1A-O3A
3	A	802	A1JZ1	OAP-CAP-CBP-CCP
3	B	802	A1JZ1	C1-C2-C3-C4
3	B	802	A1JZ1	C5B-O5B-P1A-O1A
3	B	802	A1JZ1	C3B-O3B-P3B-O7A
3	A	802	A1JZ1	O4B-C4B-C5B-O5B
3	B	802	A1JZ1	C6-C7-C8-C9
2	A	801	NAI	C2D-C1D-N1N-C6N
3	B	802	A1JZ1	C1-C2-C3-O3
3	A	802	A1JZ1	C2-C3-C4-C5
3	B	802	A1JZ1	C2-C3-C4-C5
3	A	802	A1JZ1	C7-C8-C9-C10
3	A	802	A1JZ1	C3B-C4B-C5B-O5B
3	B	802	A1JZ1	C2-C1-S1P-C2P
2	A	801	NAI	C2D-C1D-N1N-C2N
2	B	801	NAI	C2D-C1D-N1N-C6N
3	A	802	A1JZ1	CCP-O6A-P2A-O3A
3	B	802	A1JZ1	C5B-O5B-P1A-O3A
2	A	801	NAI	O4D-C1D-N1N-C6N
2	A	801	NAI	C5D-O5D-PN-O2N
3	A	802	A1JZ1	C5B-O5B-P1A-O2A
3	B	802	A1JZ1	O3-C3-C4-C5
3	A	802	A1JZ1	C5-C6-C7-C8
3	B	802	A1JZ1	C7-C8-C9-C10
2	B	801	NAI	O4D-C1D-N1N-C6N
2	A	801	NAI	O4D-C1D-N1N-C2N
3	B	802	A1JZ1	C4-C5-C6-C7
2	A	801	NAI	O4B-C4B-C5B-O5B
2	A	801	NAI	C5D-O5D-PN-O3

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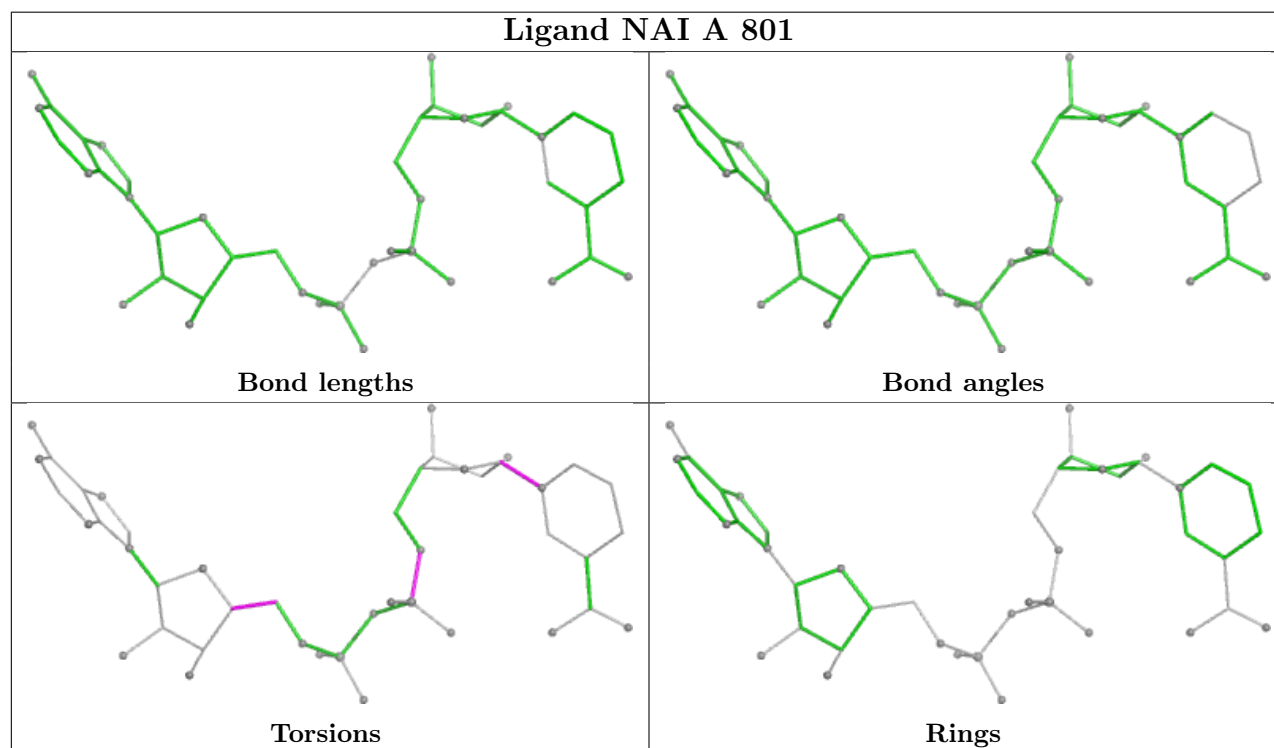
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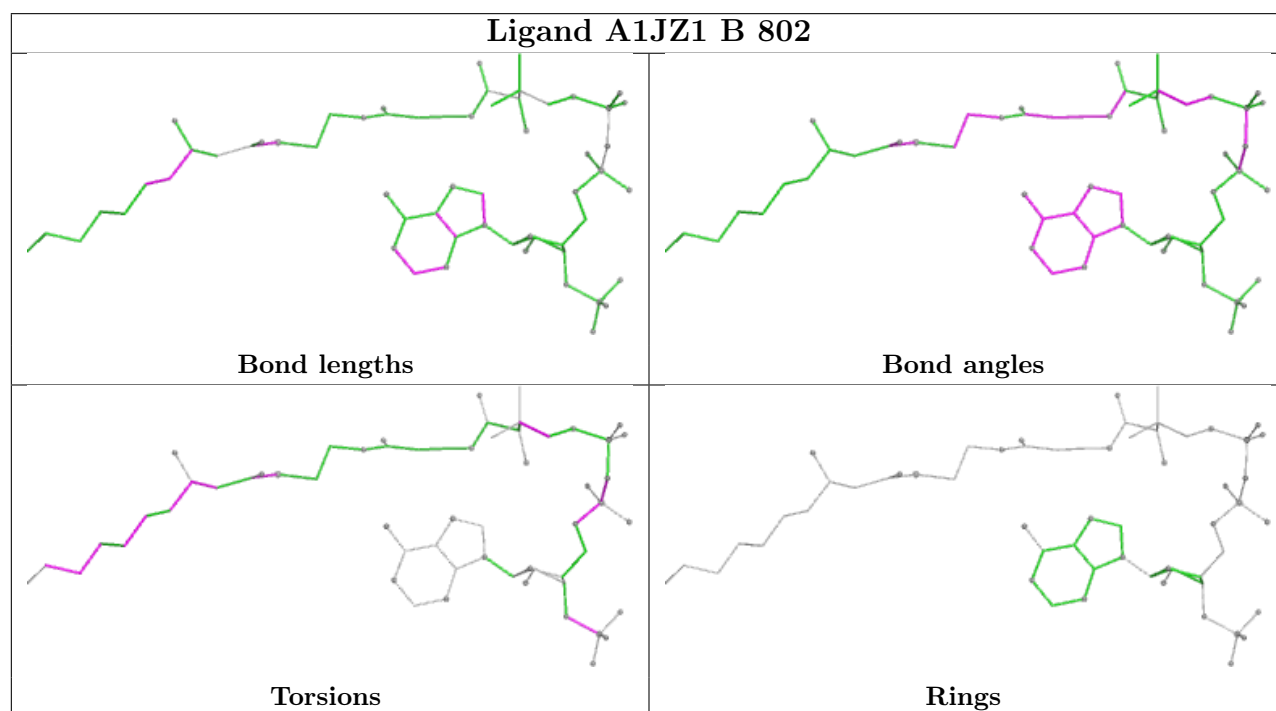
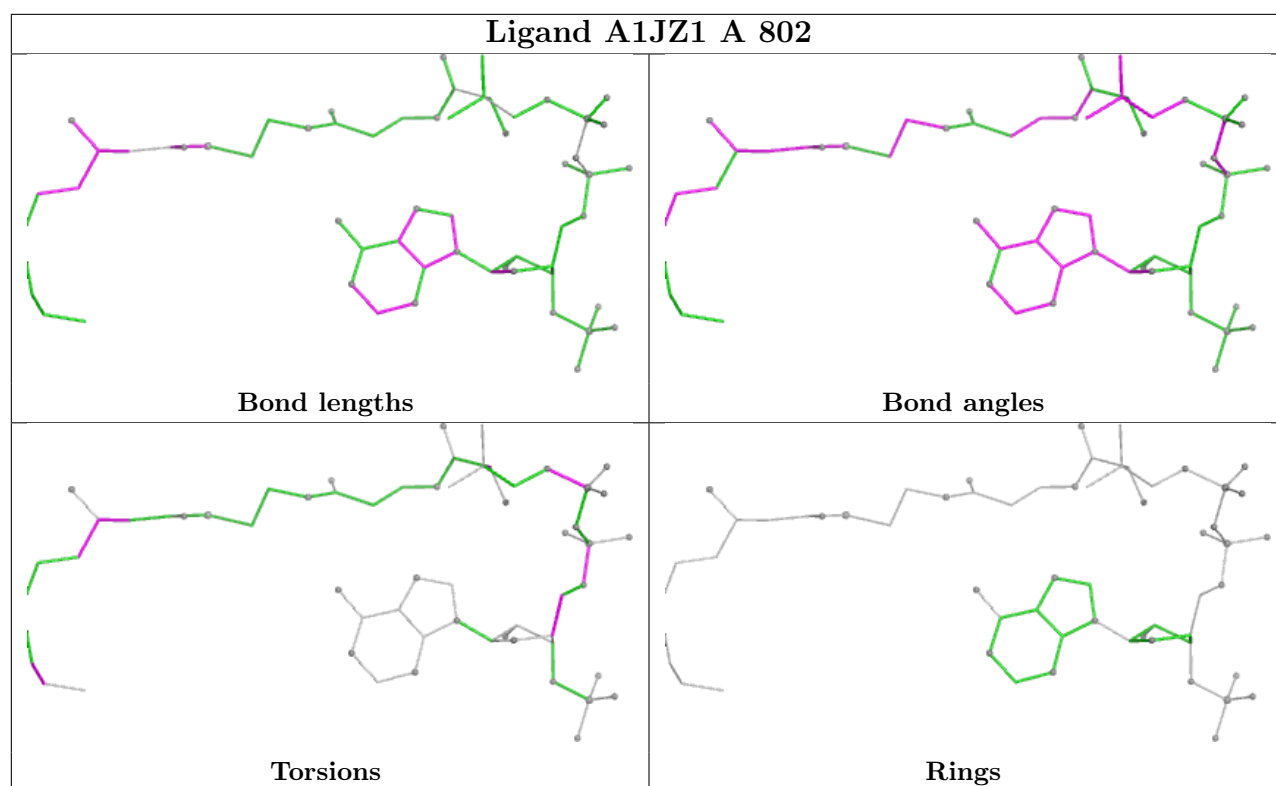
Mol	Chain	Res	Type	Atoms
2	B	801	NAI	O4B-C4B-C5B-O5B
3	B	802	A1JZ1	P2A-O3A-P1A-O1A
3	A	802	A1JZ1	CCP-O6A-P2A-O5A
3	B	802	A1JZ1	C5B-O5B-P1A-O2A
3	B	802	A1JZ1	CDP-CBP-CCP-O6A
2	B	801	NAI	O4D-C1D-N1N-C2N

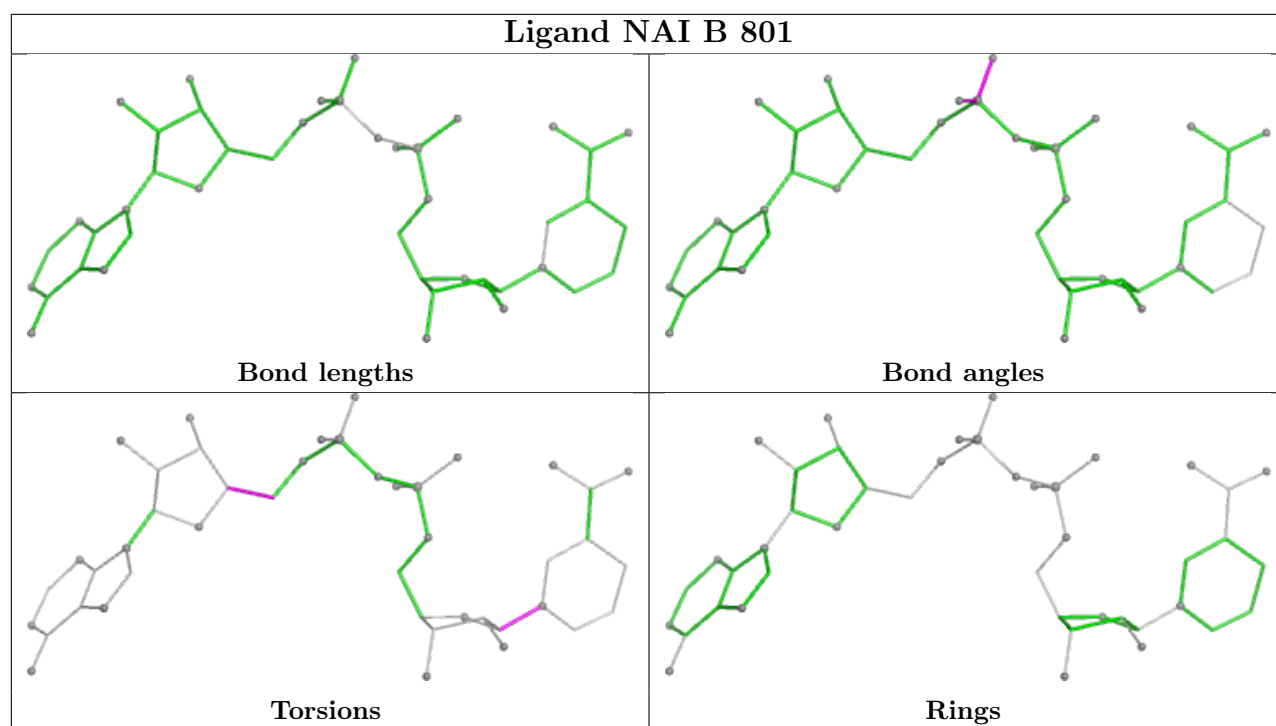
There are no ring outliers.

No monomer is involved in short contacts.

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	718/742 (96%)	-0.23	14 (1%) 66 67	37, 58, 117, 155	1 (0%)
1	B	710/742 (95%)	0.01	8 (1%) 78 79	45, 79, 120, 142	0
All	All	1428/1484 (96%)	-0.11	22 (1%) 72 73	37, 68, 118, 155	1 (0%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	69	PHE	4.9
1	B	351	HIS	3.2
1	B	73	LEU	3.2
1	A	358	ALA	3.1
1	A	71	PRO	2.9
1	A	721	LYS	2.8
1	A	398	LEU	2.7
1	B	257	TYR	2.5
1	B	359	LYS	2.4
1	A	72	GLY	2.4
1	B	1	MET	2.4
1	A	366	SER	2.3
1	A	720	SER	2.2
1	A	0	HIS	2.2
1	B	66	PHE	2.2
1	A	449	TYR	2.1
1	A	70	THR	2.1
1	A	6	ARG	2.0
1	A	699	ARG	2.0
1	A	369	LYS	2.0
1	B	531	LYS	2.0
1	B	64	HIS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

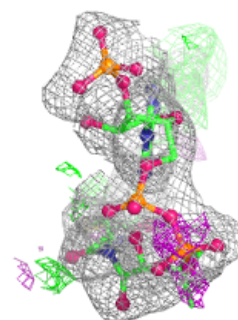
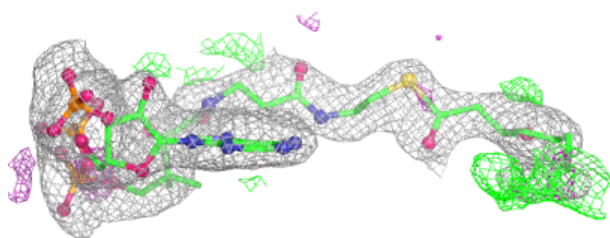
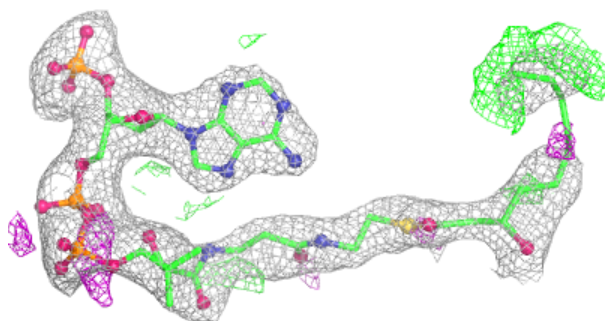
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	SO4	A	804	5/5	0.61	0.12	103,114,123,131	0
4	SO4	B	803	5/5	0.78	0.10	82,93,114,124	0
3	A1JZ1	A	802	60/60	0.85	0.14	55,89,123,130	0
2	NAI	A	801	44/44	0.90	0.11	57,73,104,119	0
3	A1JZ1	B	802	60/60	0.92	0.10	71,96,110,111	0
4	SO4	A	803	5/5	0.94	0.15	56,63,67,67	0
2	NAI	B	801	44/44	0.97	0.07	48,66,76,78	0

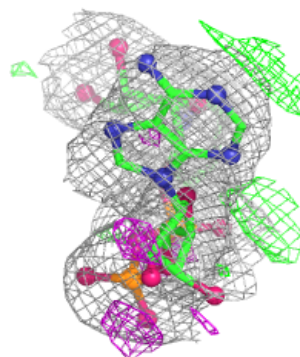
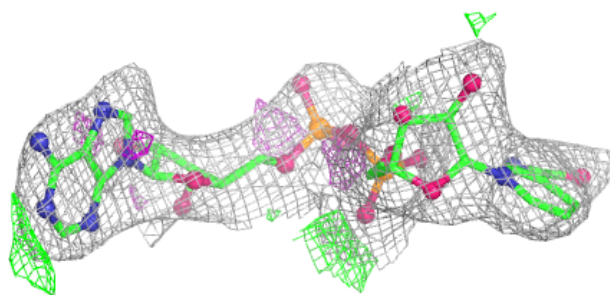
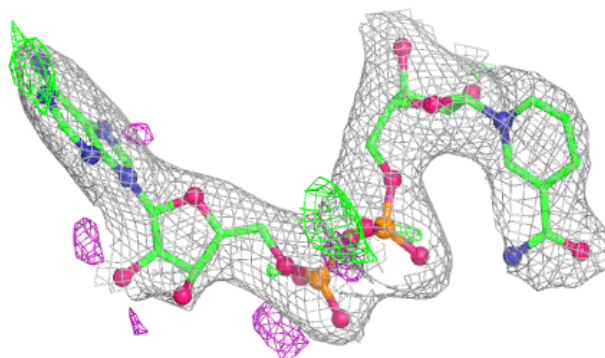
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around A1JZ1 A 802:

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

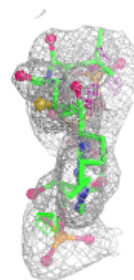
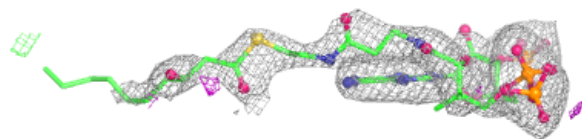
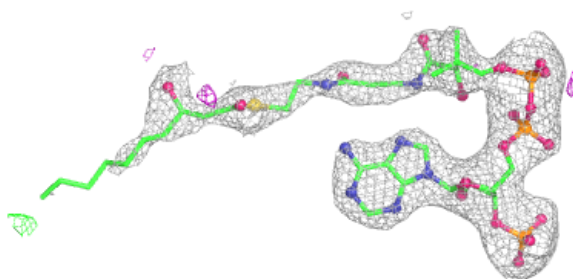
**Electron density around NAI A 801:**

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

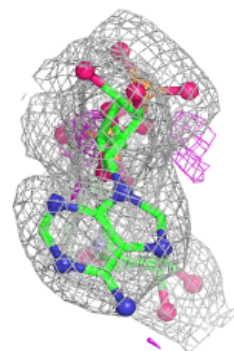
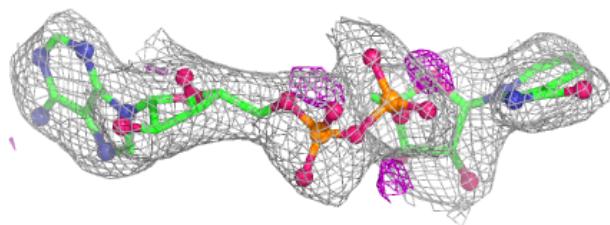
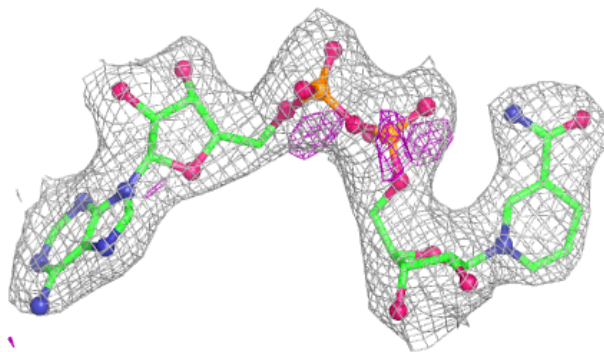


Electron density around A1JZ1 B 802:

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

**Electron density around NAI B 801:**

$2mF_o-DF_c$ (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.