



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 14, 2023 – 09:18 PM JST

PDB ID : 6AD4
Title : Rat Xanthine oxidoreductase, D428A variant, NADH bound form
Authors : Okamoto, K.; Kawaguchi, Y.
Deposited on : 2018-07-30
Resolution : 1.80 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 : 4.02b-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

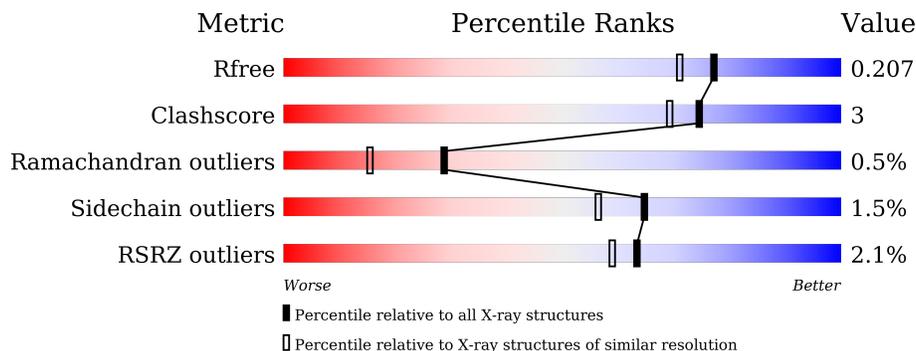
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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	1331	
1	B	1331	

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 21605 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.

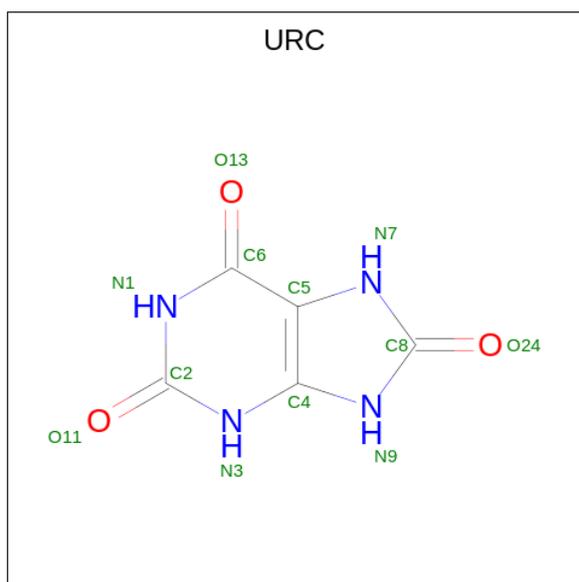
- Molecule 1 is a protein called Xanthine dehydrogenase/oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	1300	Total 10048	C 6366	N 1731	O 1886	S 65	0	0	0
1	B	1291	Total 9968	C 6319	N 1716	O 1869	S 64	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	428	ALA	ASP	see sequence details	UNP P22985
B	428	ALA	ASP	see sequence details	UNP P22985

- Molecule 2 is URIC ACID (three-letter code: URC) (formula: C₅H₄N₄O₃).



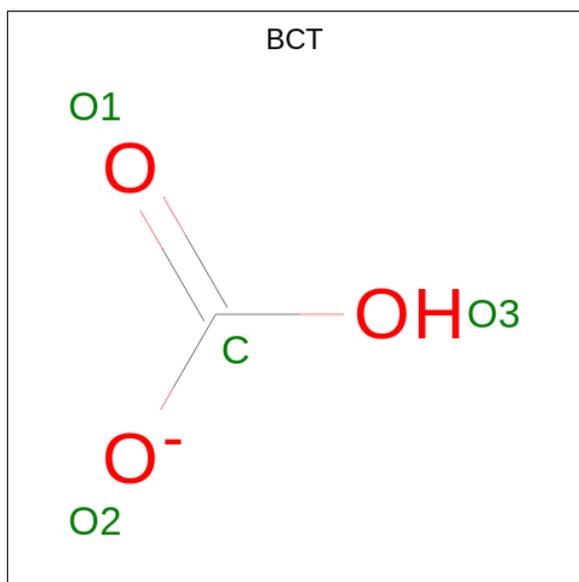
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	Total 12	C 5	N 4	O 3	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	B	1	12	5	4	3	0	0

- Molecule 3 is BICARBONATE ION (three-letter code: BCT) (formula: CHO_3).

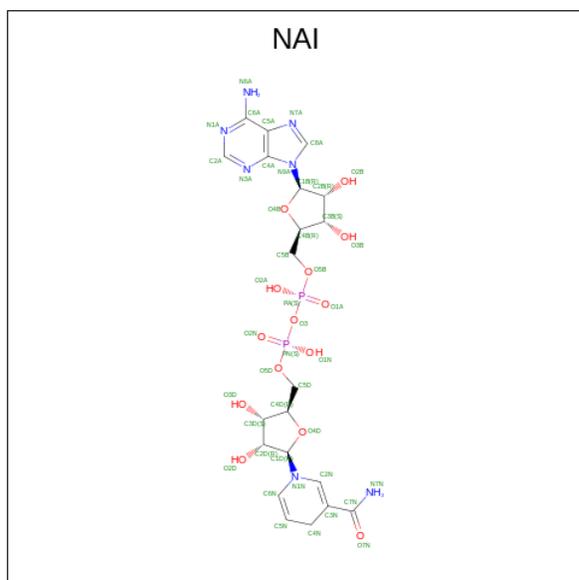


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	4	1	3	0	0
3	B	1	4	1	3	0	0

- Molecule 4 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe_2S_2).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
5	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
5	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 6 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (three-letter code: NAI) (formula: C₂₁H₂₉N₇O₁₄P₂).



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

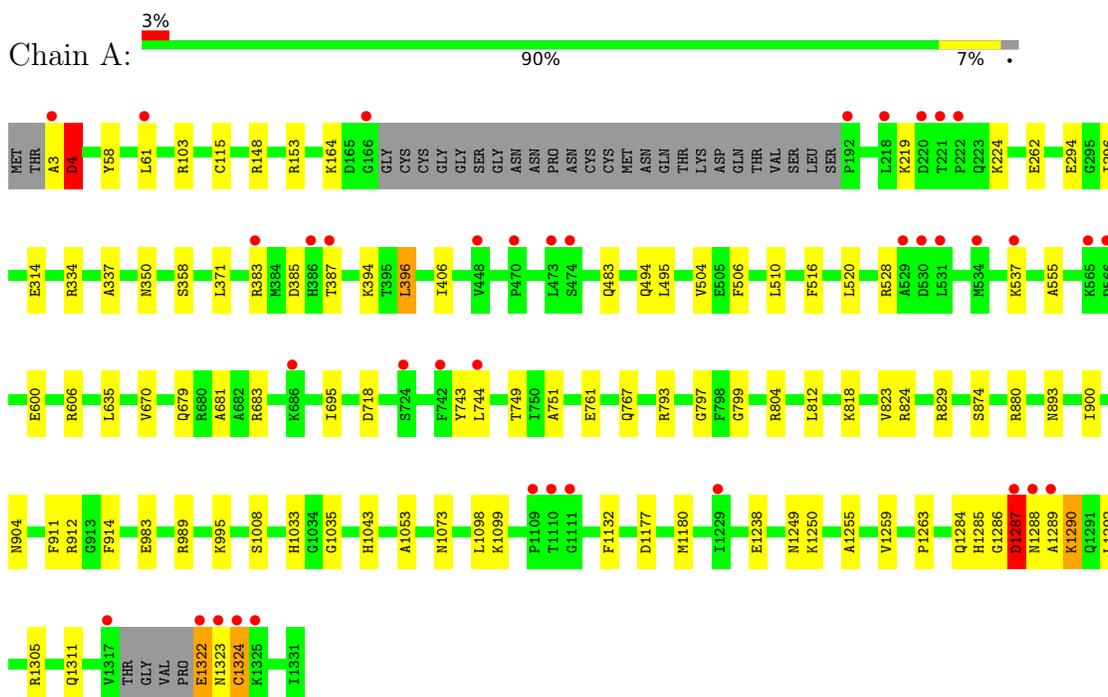
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	621	Total	O	0	0
			621	621		
7	B	726	Total	O	0	0
			726	726		

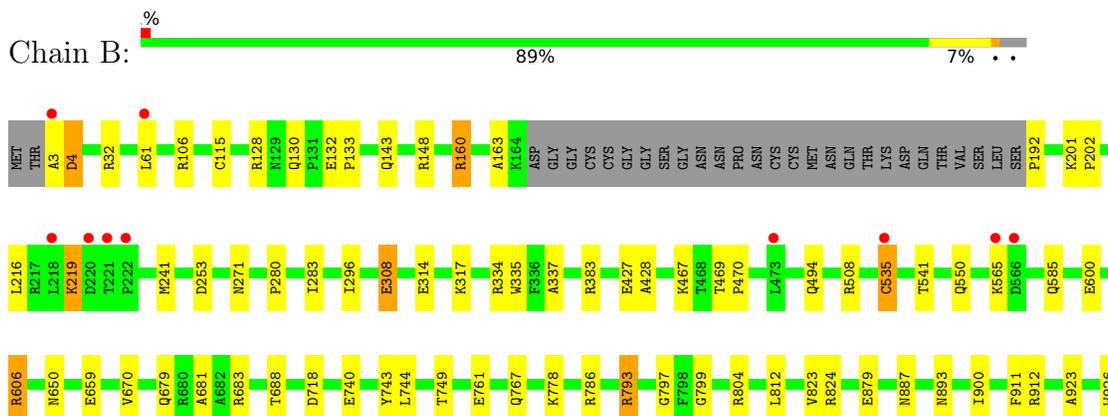
3 Residue-property plots

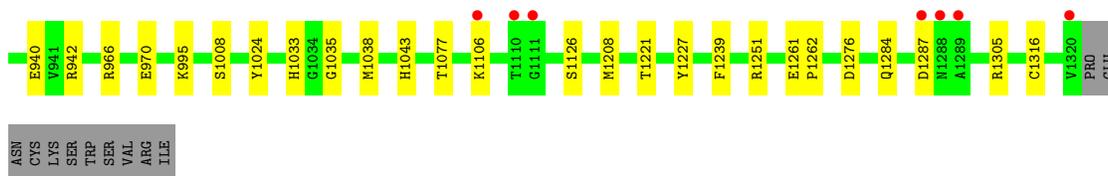
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



- Molecule 1: Xanthine dehydrogenase/oxidase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	98.72Å 138.33Å 222.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.36 – 1.80 49.36 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.36-1.80) 99.7 (49.36-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.57 (at 1.79Å)	Xtrriage
Refinement program	REFMAC 5.8.0222	Depositor
R, R_{free}	0.164 , 0.198 0.175 , 0.207	Depositor DCC
R_{free} test set	8587 reflections (3.04%)	wwPDB-VP
Wilson B-factor (Å ²)	18.9	Xtrriage
Anisotropy	0.016	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	21605	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.38% 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: BCT, FAD, NAI, FES, URC

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	0.71	4/10259 (0.0%)	0.93	7/13879 (0.1%)
1	B	0.77	12/10178 (0.1%)	0.96	12/13773 (0.1%)
All	All	0.74	16/20437 (0.1%)	0.94	19/27652 (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	6
1	B	0	11
All	All	0	17

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	600	GLU	CD-OE2	10.44	1.37	1.25
1	A	761	GLU	CD-OE1	9.42	1.36	1.25
1	B	308	GLU	CD-OE2	-8.86	1.16	1.25
1	B	740	GLU	CD-OE1	8.22	1.34	1.25
1	A	600	GLU	CD-OE2	7.90	1.34	1.25
1	A	793	ARG	CZ-NH1	7.49	1.42	1.33
1	B	761	GLU	CD-OE1	7.44	1.33	1.25
1	B	793	ARG	CZ-NH1	7.39	1.42	1.33
1	B	427	GLU	CD-OE2	7.30	1.33	1.25
1	B	940	GLU	CD-OE2	-6.71	1.18	1.25
1	A	983	GLU	CD-OE2	6.57	1.32	1.25
1	B	1276	ASP	CG-OD2	6.06	1.39	1.25
1	B	879	GLU	CD-OE1	-6.06	1.19	1.25
1	B	740	GLU	CD-OE2	5.37	1.31	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	970	GLU	CD-OE1	5.06	1.31	1.25
1	B	793	ARG	NE-CZ	5.03	1.39	1.33

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	793	ARG	NE-CZ-NH2	-40.54	100.03	120.30
1	A	793	ARG	NE-CZ-NH2	-40.41	100.09	120.30
1	B	793	ARG	NE-CZ-NH1	36.43	138.51	120.30
1	A	793	ARG	NE-CZ-NH1	36.20	138.40	120.30
1	B	793	ARG	CD-NE-CZ	12.12	140.57	123.60
1	A	793	ARG	CD-NE-CZ	11.87	140.22	123.60
1	B	160	ARG	NE-CZ-NH2	-6.79	116.91	120.30
1	B	786	ARG	NE-CZ-NH1	-6.69	116.96	120.30
1	B	541	THR	CA-CB-CG2	-6.64	103.11	112.40
1	B	160	ARG	NE-CZ-NH1	6.27	123.43	120.30
1	B	942	ARG	NE-CZ-NH2	-5.82	117.39	120.30
1	B	606	ARG	NE-CZ-NH1	-5.80	117.40	120.30
1	A	989	ARG	NE-CZ-NH2	-5.64	117.48	120.30
1	B	585	GLN	CA-C-N	-5.32	105.50	117.20
1	B	966	ARG	NE-CZ-NH2	-5.29	117.65	120.30
1	A	829	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	B	106	ARG	NE-CZ-NH2	-5.19	117.71	120.30
1	A	334	ARG	NE-CZ-NH2	-5.14	117.73	120.30
1	A	103	ARG	NE-CZ-NH2	-5.01	117.79	120.30

There are no chirality outliers.

All (17) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1287	ASP	Peptide
1	A	1305	ARG	Sidechain
1	A	148	ARG	Sidechain
1	A	3	ALA	Peptide
1	A	528	ARG	Sidechain
1	A	804	ARG	Sidechain
1	B	1251	ARG	Sidechain
1	B	128	ARG	Sidechain
1	B	1305	ARG	Sidechain
1	B	148	ARG	Sidechain
1	B	160	ARG	Sidechain
1	B	3	ALA	Peptide

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Mol	Chain	Res	Type	Group
1	B	32	ARG	Sidechain
1	B	334	ARG	Sidechain
1	B	383	ARG	Sidechain
1	B	535	CYS	Peptide
1	B	804	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	10048	0	10053	80	0
1	B	9968	0	9982	53	0
2	A	12	0	4	0	0
2	B	12	0	4	0	0
3	A	4	0	0	0	0
3	B	4	0	0	0	0
4	A	8	0	0	0	0
4	B	8	0	0	1	0
5	A	53	0	31	1	0
5	B	53	0	31	1	0
6	A	44	0	27	0	0
6	B	44	0	27	0	0
7	A	621	0	0	6	0
7	B	726	0	0	4	0
All	All	21605	0	20159	131	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 (131) 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:1180:MET:SD	7:A:2150:HOH:O	1.96	1.18
1:A:385:ASP:OD1	1:A:387:THR:HG22	1.44	1.15
1:B:718:ASP:H	1:B:893:ASN:HD22	1.18	0.91
1:B:749:THR:HG22	1:B:812:LEU:HD23	1.52	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:271:ASN:ND2	1:B:683:ARG:HH11	1.76	0.83
1:A:695:ILE:H	1:A:904:ASN:HD22	1.26	0.82
1:A:751:ALA:CB	1:A:812:LEU:CD2	2.59	0.81
1:A:751:ALA:N	1:A:812:LEU:HD21	1.97	0.80
1:A:751:ALA:CB	1:A:812:LEU:HD23	2.13	0.79
1:A:1323:ASN:HA	1:A:1324:CYS:CB	2.13	0.78
1:A:1322:GLU:HB2	1:A:1323:ASN:HB2	1.67	0.77
1:B:216:LEU:O	1:B:219:LYS:HG3	1.87	0.74
1:A:1289:ALA:O	1:A:1290:LYS:HB3	1.88	0.74
1:A:483:GLN:HE22	1:A:1311:GLN:HE22	1.37	0.73
1:A:718:ASP:H	1:A:893:ASN:HD22	1.38	0.72
1:A:1033:HIS:HE1	1:A:1043:HIS:HD2	1.35	0.72
1:B:130:GLN:HE21	1:B:132:GLU:H	1.37	0.71
1:A:751:ALA:HB2	1:A:812:LEU:CD2	2.21	0.70
1:A:751:ALA:H	1:A:812:LEU:HD21	1.54	0.70
1:B:241:MET:HE2	1:B:283:ILE:HG21	1.74	0.70
1:A:1323:ASN:CA	1:A:1324:CYS:HB2	2.22	0.69
1:B:812:LEU:CD1	1:B:823:VAL:HG12	2.22	0.69
1:A:749:THR:HG22	1:A:812:LEU:HD22	1.75	0.68
1:A:749:THR:O	1:A:812:LEU:HD13	1.93	0.68
1:B:995:LYS:NZ	1:B:1284:GLN:HE21	1.91	0.68
1:B:192:PRO:HD2	7:B:3620:HOH:O	1.94	0.68
1:A:262:GLU:OE1	7:A:1601:HOH:O	2.11	0.67
1:B:718:ASP:H	1:B:893:ASN:ND2	1.92	0.67
1:B:749:THR:HG22	1:B:812:LEU:CD2	2.22	0.67
1:A:751:ALA:HB3	1:A:812:LEU:HD23	1.77	0.67
1:A:606:ARG:HD3	1:A:679:GLN:HA	1.78	0.66
1:A:751:ALA:HB2	1:A:812:LEU:HD23	1.76	0.66
1:A:1323:ASN:CA	1:A:1324:CYS:CB	2.73	0.66
1:B:688:THR:HG23	7:B:3581:HOH:O	1.95	0.65
1:B:812:LEU:HD11	1:B:823:VAL:HG12	1.77	0.65
1:A:1323:ASN:HA	1:A:1324:CYS:HB3	1.78	0.65
1:B:606:ARG:HD3	1:B:679:GLN:HA	1.79	0.65
1:A:1323:ASN:HA	1:A:1324:CYS:HB2	1.80	0.64
1:B:749:THR:O	1:B:812:LEU:HD22	1.98	0.64
1:B:253:ASP:HB2	7:B:3312:HOH:O	1.98	0.64
1:A:995:LYS:NZ	1:A:1284:GLN:HE21	1.96	0.63
1:A:1033:HIS:CE1	1:A:1043:HIS:HD2	2.15	0.62
1:A:812:LEU:HD11	1:A:824:ARG:HA	1.79	0.62
1:A:767:GLN:HE22	1:A:799:GLY:H	1.48	0.62
1:B:812:LEU:HD21	1:B:823:VAL:O	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:767:GLN:HE22	1:B:799:GLY:H	1.46	0.61
1:A:1033:HIS:HD2	1:A:1035:GLY:H	1.47	0.61
1:B:812:LEU:HD11	1:B:823:VAL:C	2.21	0.61
1:A:751:ALA:HB3	1:A:812:LEU:CD2	2.30	0.61
1:A:1323:ASN:N	1:A:1324:CYS:HB2	2.15	0.61
1:A:695:ILE:H	1:A:904:ASN:ND2	1.99	0.60
1:B:115:CYS:HB3	1:B:744:LEU:HD22	1.82	0.60
1:A:1033:HIS:CE1	1:A:1043:HIS:CD2	2.92	0.58
1:A:1285:HIS:HD2	7:A:2177:HOH:O	1.87	0.58
1:A:1033:HIS:HE1	1:A:1043:HIS:CD2	2.19	0.57
1:A:718:ASP:H	1:A:893:ASN:ND2	2.01	0.57
1:B:271:ASN:ND2	1:B:683:ARG:NH1	2.51	0.57
1:B:115:CYS:HB3	1:B:744:LEU:CD2	2.34	0.56
1:B:995:LYS:HZ1	1:B:1284:GLN:HE21	1.54	0.56
1:A:1322:GLU:CB	1:A:1323:ASN:HB2	2.34	0.55
1:B:1033:HIS:HD1	1:B:1035:GLY:H	1.54	0.54
1:A:115:CYS:HB3	1:A:744:LEU:HD22	1.89	0.54
1:B:143:GLN:HE21	1:B:428:ALA:H	1.55	0.54
1:A:1033:HIS:CD2	1:A:1035:GLY:H	2.26	0.52
1:A:1073:ASN:HD21	1:B:1024:TYR:HA	1.75	0.52
1:B:143:GLN:HE22	1:B:335:TRP:HA	1.74	0.51
1:A:4:ASP:HB2	7:A:1744:HOH:O	2.10	0.50
1:A:1289:ALA:O	1:A:1290:LYS:CB	2.59	0.50
1:A:58:TYR:CE1	1:A:219:LYS:HD2	2.47	0.50
1:A:812:LEU:HD11	1:A:824:ARG:CA	2.42	0.49
1:A:812:LEU:HD11	1:A:823:VAL:O	2.12	0.49
1:A:555:ALA:O	1:A:1238:GLU:HA	2.13	0.49
1:A:874:SER:HB3	1:A:900:ILE:HG21	1.95	0.49
1:A:749:THR:O	1:A:812:LEU:CD1	2.59	0.48
1:A:1053:ALA:O	1:A:1098:LEU:HD11	2.13	0.48
1:B:812:LEU:HD12	1:B:823:VAL:HG12	1.93	0.48
1:A:751:ALA:CB	1:A:812:LEU:HD21	2.40	0.48
1:A:58:TYR:CZ	1:A:219:LYS:HD2	2.49	0.47
1:A:371:LEU:HD21	1:A:406:ILE:HD12	1.97	0.47
1:A:995:LYS:HZ3	1:A:1284:GLN:HE21	1.62	0.47
1:B:296:ILE:CD1	1:B:314:GLU:HG3	2.45	0.47
1:A:350:ASN:ND2	1:A:358:SER:OG	2.45	0.47
1:A:670:VAL:HG11	1:A:681:ALA:HB3	1.96	0.47
1:A:1287:ASP:O	1:A:1287:ASP:CG	2.53	0.46
1:B:1033:HIS:NE2	1:B:1043:HIS:HD2	2.13	0.46
1:A:337:ALA:HB2	5:A:1505:FAD:C6	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:508:ARG:NH1	1:B:1316:CYS:O	2.40	0.46
1:B:744:LEU:HD22	4:B:3001:FES:S1	2.56	0.46
1:B:812:LEU:HD11	1:B:824:ARG:N	2.31	0.45
1:B:1221:THR:HG22	1:B:1227:TYR:HB2	1.98	0.45
1:B:670:VAL:HG11	1:B:681:ALA:HB3	1.98	0.45
1:A:1292:LEU:HA	7:A:1914:HOH:O	2.17	0.45
1:A:516:PHE:CZ	1:A:520:LEU:HD11	2.53	0.44
1:B:767:GLN:HE22	1:B:799:GLY:N	2.13	0.44
1:B:900:ILE:HD12	1:B:900:ILE:N	2.33	0.44
1:A:115:CYS:HB3	1:A:744:LEU:CD2	2.48	0.43
1:A:1180:MET:HE1	1:A:1263:PRO:HB3	2.00	0.43
1:B:469:THR:N	1:B:470:PRO:CD	2.82	0.43
1:B:1208:MET:HE3	1:B:1208:MET:HB3	1.90	0.43
1:A:495:LEU:HB2	1:A:504:VAL:HG13	1.99	0.43
1:A:1180:MET:CE	1:A:1263:PRO:HB3	2.49	0.43
1:A:880:ARG:HD2	1:A:914:PHE:HB3	2.01	0.43
1:B:1261:GLU:N	1:B:1262:PRO:CD	2.82	0.43
1:A:635:LEU:HD21	1:A:818:LYS:HD3	2.01	0.42
1:B:995:LYS:HZ3	1:B:1284:GLN:HE21	1.61	0.42
1:A:995:LYS:HZ1	1:A:1284:GLN:HE21	1.66	0.42
1:A:1259:VAL:HG22	1:A:1259:VAL:O	2.20	0.42
1:B:337:ALA:HB2	5:B:3005:FAD:C6	2.49	0.42
1:A:751:ALA:N	1:A:812:LEU:CD2	2.76	0.42
1:B:133:PRO:O	1:B:163:ALA:HA	2.19	0.42
1:B:271:ASN:HD22	1:B:683:ARG:NH1	2.18	0.42
1:A:1043:HIS:HE1	7:A:1870:HOH:O	2.03	0.41
1:A:1249:ASN:O	1:A:1255:ALA:HA	2.21	0.41
1:A:153:ARG:HD2	1:A:153:ARG:C	2.40	0.41
1:A:296:ILE:CD1	1:A:314:GLU:HG3	2.50	0.41
1:B:1033:HIS:NE2	1:B:1043:HIS:CD2	2.88	0.41
1:B:201:LYS:HA	1:B:202:PRO:HD3	1.90	0.41
1:B:793:ARG:NH1	7:B:3116:HOH:O	2.53	0.41
1:A:296:ILE:HD11	1:A:314:GLU:HG3	2.03	0.41
1:A:506:PHE:CZ	1:A:510:LEU:HD11	2.56	0.41
1:A:718:ASP:N	1:A:893:ASN:HD22	2.13	0.41
1:A:767:GLN:HE22	1:A:799:GLY:N	2.15	0.41
1:B:650:ASN:OD1	1:B:778:LYS:NZ	2.54	0.41
1:A:1132:PHE:CG	1:B:1126:SER:HB2	2.57	0.40
1:B:130:GLN:NE2	1:B:132:GLU:H	2.13	0.40
1:A:679:GLN:HG2	1:A:683:ARG:HH11	1.86	0.40
1:B:1038:MET:CE	1:B:1077:THR:CG2	2.99	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:749:THR:CG2	1:A:812:LEU:HD22	2.47	0.40
1:A:396:LEU:C	1:A:396:LEU:HD12	2.42	0.40
1:B:923:ALA:HA	1:B:926:TRP:NE1	2.37	0.40
1:B:565:LYS:HD3	1:B:565:LYS:HA	2.01	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	1294/1331 (97%)	1245 (96%)	41 (3%)	8 (1%)	25	12
1	B	1287/1331 (97%)	1250 (97%)	31 (2%)	6 (0%)	29	15
All	All	2581/2662 (97%)	2495 (97%)	72 (3%)	14 (0%)	29	15

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	ASP
1	A	1008	SER
1	A	1324	CYS
1	B	4	ASP
1	B	1008	SER
1	A	1288	ASN
1	A	1290	LYS
1	B	912	ARG
1	A	912	ARG
1	A	797	GLY
1	B	797	GLY
1	B	887	ASN
1	B	1287	ASP
1	A	1286	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	1097/1123 (98%)	1080 (98%)	17 (2%)	65	56
1	B	1088/1123 (97%)	1073 (99%)	15 (1%)	67	59
All	All	2185/2246 (97%)	2153 (98%)	32 (2%)	65	56

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ASP
1	A	61	LEU
1	A	164	LYS
1	A	224	LYS
1	A	294	GLU
1	A	383	ARG
1	A	394	LYS
1	A	396	LEU
1	A	494	GLN
1	A	537	LYS
1	A	743	TYR
1	A	911	PHE
1	A	1099	LYS
1	A	1177	ASP
1	A	1250	LYS
1	A	1287	ASP
1	A	1322	GLU
1	B	4	ASP
1	B	61	LEU
1	B	219	LYS
1	B	280	PRO
1	B	308	GLU
1	B	317	LYS
1	B	467	LYS
1	B	494	GLN
1	B	535	CYS
1	B	550	GLN

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Mol	Chain	Res	Type
1	B	659	GLU
1	B	743	TYR
1	B	911	PHE
1	B	1106	LYS
1	B	1239	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	ASN
1	A	350	ASN
1	A	585	GLN
1	A	767	GLN
1	A	893	ASN
1	A	904	ASN
1	A	956	ASN
1	A	1016	GLN
1	A	1033	HIS
1	A	1043	HIS
1	A	1073	ASN
1	A	1284	GLN
1	A	1285	HIS
1	A	1311	GLN
1	B	130	GLN
1	B	143	GLN
1	B	145	ASN
1	B	271	ASN
1	B	350	ASN
1	B	550	GLN
1	B	585	GLN
1	B	767	GLN
1	B	893	ASN
1	B	956	ASN
1	B	1016	GLN
1	B	1043	HIS
1	B	1284	GLN

5.3.3 RNA

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](#)

12 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	BCT	A	1502	-	2,3,3	0.94	0	2,3,3	0.25	0
4	FES	A	1503	1	0,4,4	-	-	-	-	-
5	FAD	B	3005	-	53,58,58	1.49	8 (15%)	68,89,89	1.33	13 (19%)
6	NAI	A	1506	-	42,48,48	1.29	5 (11%)	47,73,73	1.40	7 (14%)
3	BCT	B	3004	-	2,3,3	1.03	0	2,3,3	2.36	1 (50%)
2	URC	A	1501	-	8,13,13	3.28	5 (62%)	5,19,19	6.69	4 (80%)
2	URC	B	3003	-	8,13,13	2.92	5 (62%)	5,19,19	6.32	4 (80%)
6	NAI	B	3006	-	42,48,48	1.33	6 (14%)	47,73,73	1.28	5 (10%)
4	FES	A	1504	1	0,4,4	-	-	-	-	-
5	FAD	A	1505	-	53,58,58	1.56	12 (22%)	68,89,89	1.23	6 (8%)
4	FES	B	3002	1	0,4,4	-	-	-	-	-
4	FES	B	3001	1	0,4,4	-	-	-	-	-

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	FAD	B	3005	-	-	2/30/50/50	0/6/6/6

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	FES	A	1503	1	-	-	0/1/1/1
6	NAI	A	1506	-	-	3/25/72/72	0/5/5/5
2	URC	A	1501	-	-	-	0/2/2/2
2	URC	B	3003	-	-	-	0/2/2/2
6	NAI	B	3006	-	-	4/25/72/72	0/5/5/5
4	FES	A	1504	1	-	-	0/1/1/1
5	FAD	A	1505	-	-	5/30/50/50	0/6/6/6
4	FES	B	3002	1	-	-	0/1/1/1
4	FES	B	3001	1	-	-	0/1/1/1

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1501	URC	C5-C6	5.98	1.51	1.41
2	B	3003	URC	C5-C6	4.51	1.49	1.41
5	B	3005	FAD	C4X-N5	4.37	1.39	1.30
5	B	3005	FAD	O4B-C1B	4.00	1.46	1.41
2	A	1501	URC	C5-C4	3.96	1.51	1.40
2	B	3003	URC	C5-C4	3.91	1.51	1.40
5	B	3005	FAD	C9A-C5X	3.81	1.47	1.41
2	B	3003	URC	C4-N9	3.80	1.41	1.34
6	A	1506	NAI	C6N-C5N	3.68	1.39	1.33
5	B	3005	FAD	C2B-C1B	-3.40	1.48	1.53
2	A	1501	URC	C4-N9	3.39	1.41	1.34
5	A	1505	FAD	C2B-C1B	-3.28	1.48	1.53
6	B	3006	NAI	O4B-C1B	3.12	1.45	1.41
2	A	1501	URC	O13-C6	3.08	1.32	1.24
6	A	1506	NAI	C2N-C3N	3.04	1.43	1.34
5	A	1505	FAD	C9A-C5X	3.00	1.46	1.41
2	B	3003	URC	C6-N1	2.97	1.38	1.33
5	A	1505	FAD	C4X-N5	2.95	1.36	1.30
5	B	3005	FAD	C8-C7	2.91	1.48	1.40
5	A	1505	FAD	O4B-C1B	2.86	1.45	1.41
6	B	3006	NAI	C2B-C1B	-2.75	1.49	1.53
6	B	3006	NAI	C5A-C4A	2.69	1.48	1.40
5	A	1505	FAD	O2-C2	2.68	1.29	1.24
2	A	1501	URC	C6-N1	2.68	1.37	1.33
6	B	3006	NAI	C7N-C3N	2.63	1.54	1.48
5	A	1505	FAD	C5X-N5	-2.62	1.34	1.39
5	B	3005	FAD	C10-N1	2.59	1.38	1.33
6	A	1506	NAI	C5A-C4A	2.58	1.47	1.40
5	A	1505	FAD	C8-C7	2.57	1.47	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	1506	NAI	O4B-C1B	2.52	1.44	1.41
5	A	1505	FAD	C1'-C2'	2.44	1.56	1.52
6	B	3006	NAI	C2N-C3N	2.35	1.41	1.34
5	A	1505	FAD	C7M-C7	2.30	1.55	1.51
2	B	3003	URC	O13-C6	2.25	1.30	1.24
5	A	1505	FAD	O2B-C2B	2.18	1.48	1.43
5	B	3005	FAD	O3'-C3'	2.13	1.48	1.43
5	A	1505	FAD	O3'-C3'	2.10	1.47	1.43
5	B	3005	FAD	C5X-N5	-2.06	1.35	1.39
6	B	3006	NAI	O4D-C1D	2.05	1.46	1.42
5	A	1505	FAD	O3B-C3B	2.02	1.47	1.43
6	A	1506	NAI	C2B-C1B	-2.00	1.50	1.53

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1501	URC	C4-N9-C8	9.74	107.03	101.12
2	B	3003	URC	C4-N9-C8	8.85	106.49	101.12
2	A	1501	URC	C2-N1-C6	8.70	122.49	115.14
2	B	3003	URC	C2-N1-C6	8.59	122.39	115.14
2	B	3003	URC	C5-C6-N1	-6.06	115.14	123.43
2	A	1501	URC	C5-C6-N1	-5.98	115.25	123.43
6	A	1506	NAI	N3A-C2A-N1A	-4.51	121.63	128.68
2	A	1501	URC	C4-C5-N7	-4.09	105.32	109.47
5	A	1505	FAD	O4'-C4'-C3'	-3.45	100.72	109.10
5	A	1505	FAD	C1B-N9A-C4A	-3.44	120.60	126.64
2	B	3003	URC	C4-C5-N7	-3.29	106.13	109.47
6	B	3006	NAI	N3A-C2A-N1A	-3.26	123.58	128.68
3	B	3004	BCT	O2-C-O1	3.19	127.83	119.55
5	B	3005	FAD	O4-C4-C4X	-2.86	119.02	126.60
5	B	3005	FAD	C1B-N9A-C4A	-2.84	121.64	126.64
6	B	3006	NAI	C3D-C2D-C1D	2.84	106.82	101.43
5	B	3005	FAD	C4-N3-C2	-2.82	120.43	125.64
6	A	1506	NAI	C3N-C2N-N1N	-2.80	119.10	123.10
5	B	3005	FAD	N6A-C6A-N1A	2.64	124.05	118.57
6	A	1506	NAI	PN-O3-PA	-2.64	123.77	132.83
5	A	1505	FAD	C2B-C3B-C4B	2.59	107.67	102.64
5	B	3005	FAD	O4'-C4'-C3'	-2.57	102.86	109.10
5	B	3005	FAD	N3-C2-N1	2.50	124.28	119.38
6	B	3006	NAI	C4A-C5A-N7A	-2.49	106.80	109.40
6	B	3006	NAI	PN-O3-PA	-2.49	124.27	132.83
6	A	1506	NAI	C5A-C6A-N6A	-2.48	116.59	120.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	3005	FAD	C4X-C10-N1	-2.45	119.05	124.73
6	A	1506	NAI	N6A-C6A-N1A	2.45	123.66	118.57
5	B	3005	FAD	C5'-C4'-C3'	2.42	116.88	112.20
5	A	1505	FAD	N6A-C6A-N1A	2.36	123.47	118.57
5	A	1505	FAD	C7M-C7-C8	2.29	125.44	120.74
6	B	3006	NAI	O7N-C7N-C3N	-2.27	116.62	120.90
5	B	3005	FAD	C4-C4X-N5	2.24	121.42	118.23
5	B	3005	FAD	C4X-C4-N3	2.20	118.78	113.19
6	A	1506	NAI	O2B-C2B-C1B	-2.19	102.78	110.85
5	B	3005	FAD	C10-C4X-N5	-2.08	120.44	124.86
5	A	1505	FAD	C9A-N10-C10	-2.06	117.56	120.77
5	B	3005	FAD	C4X-C10-N10	2.02	119.44	116.48
5	B	3005	FAD	O2-C2-N1	-2.01	118.49	121.83
6	A	1506	NAI	O5D-C5D-C4D	-2.01	102.09	108.99

There are no chirality outliers.

All (14) torsion outliers are listed below:

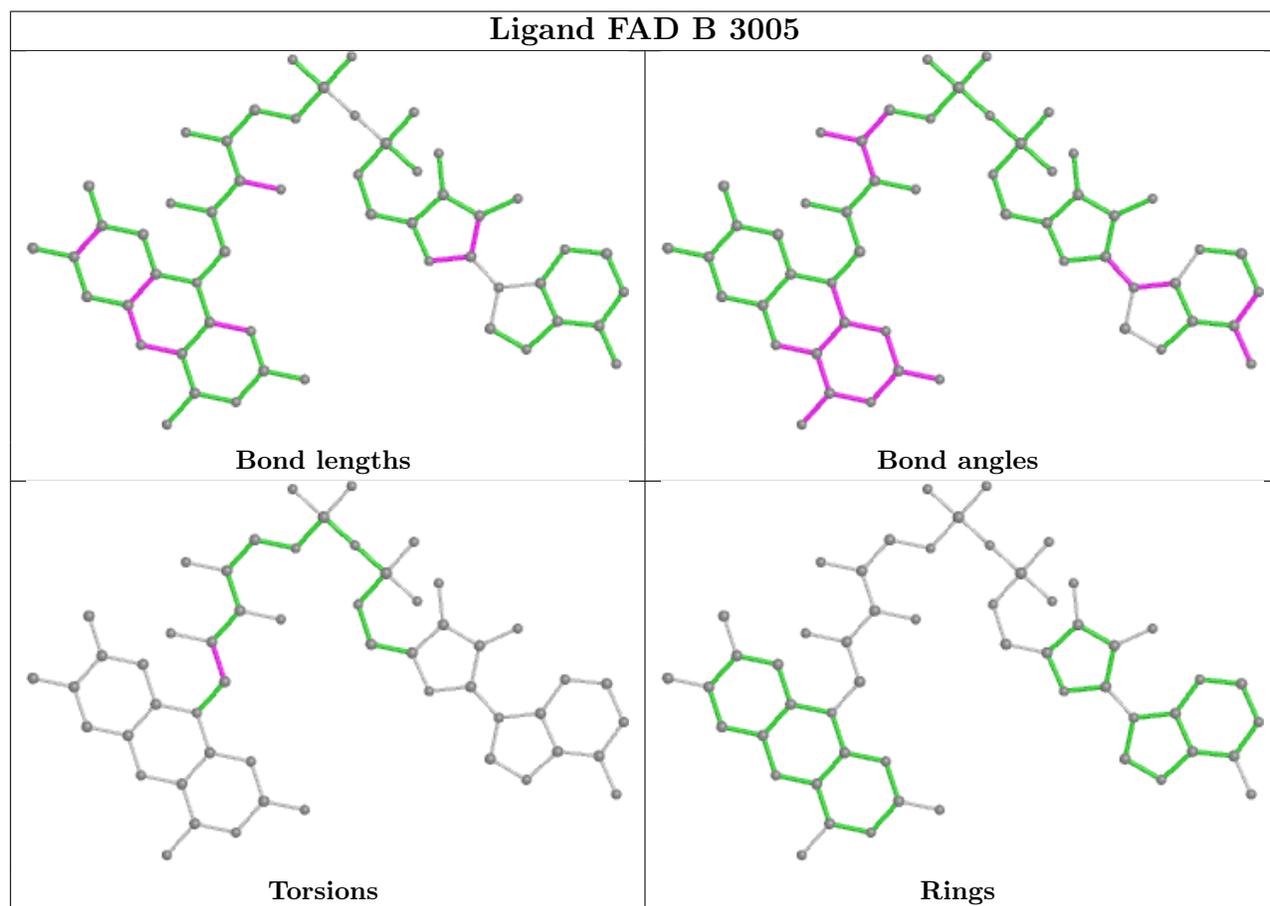
Mol	Chain	Res	Type	Atoms
5	A	1505	FAD	N10-C1'-C2'-O2'
5	A	1505	FAD	N10-C1'-C2'-C3'
5	B	3005	FAD	N10-C1'-C2'-O2'
5	B	3005	FAD	N10-C1'-C2'-C3'
5	A	1505	FAD	C2'-C3'-C4'-O4'
5	A	1505	FAD	C2'-C3'-C4'-C5'
6	A	1506	NAI	C2D-C1D-N1N-C6N
6	A	1506	NAI	O4D-C1D-N1N-C6N
6	B	3006	NAI	C2D-C1D-N1N-C6N
6	B	3006	NAI	O4D-C1D-N1N-C6N
6	B	3006	NAI	PA-O3-PN-O1N
6	A	1506	NAI	O4D-C4D-C5D-O5D
6	B	3006	NAI	O4D-C4D-C5D-O5D
5	A	1505	FAD	C2'-C1'-N10-C10

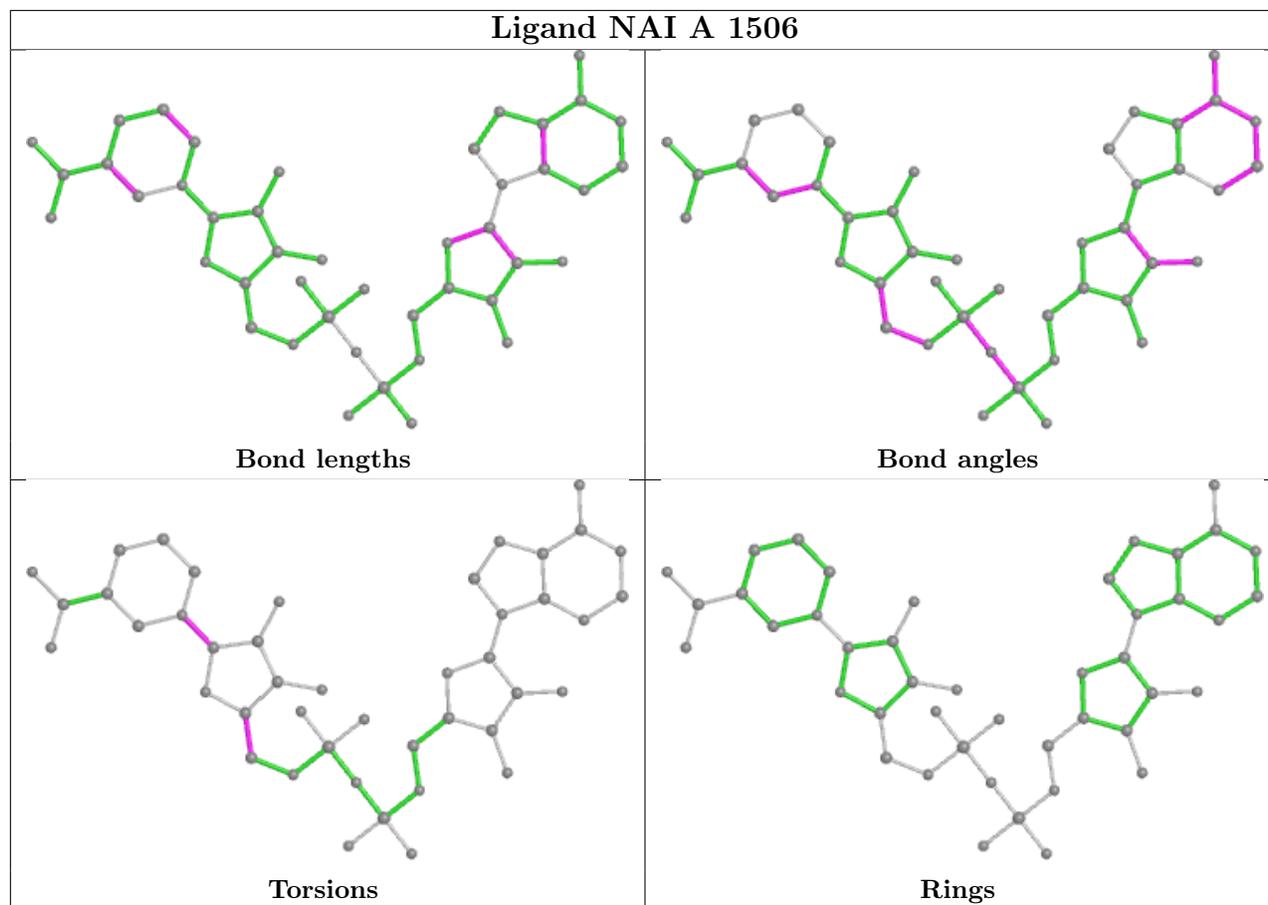
There are no ring outliers.

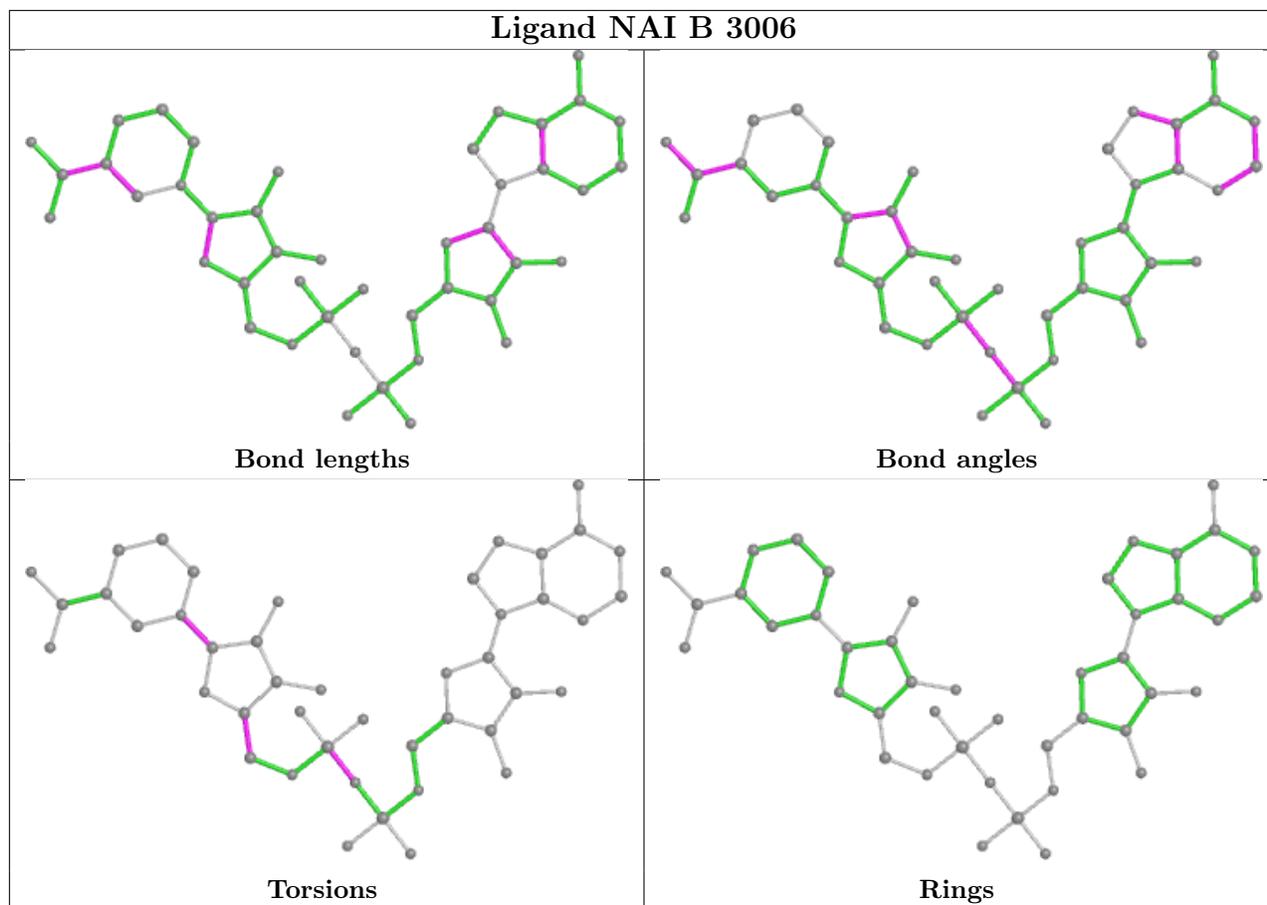
3 monomers are involved in 3 short contacts:

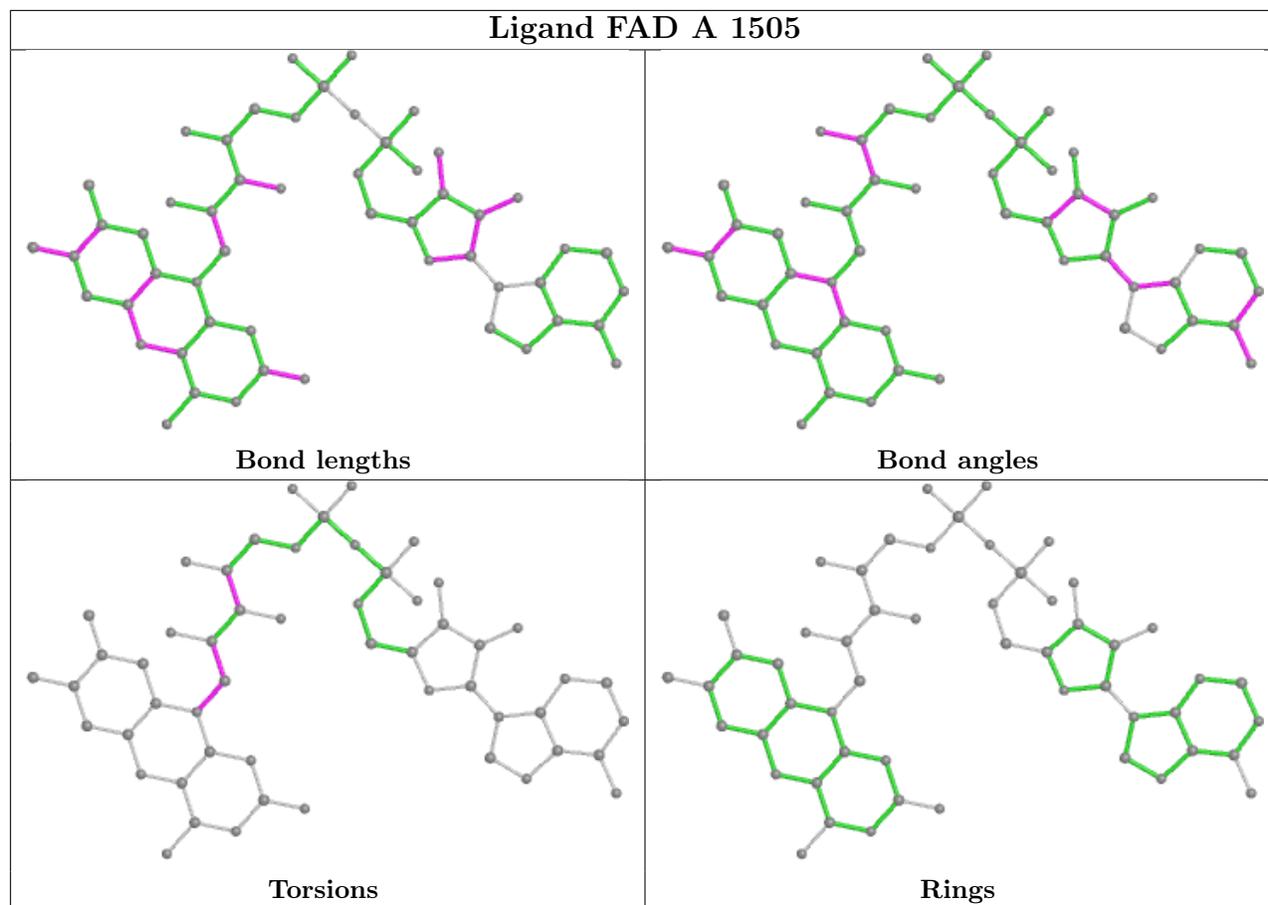
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	3005	FAD	1	0
5	A	1505	FAD	1	0
4	B	3001	FES	1	0

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	1300/1331 (97%)	0.05	38 (2%) 51 46	11, 21, 43, 81	0
1	B	1291/1331 (96%)	-0.14	17 (1%) 77 74	11, 17, 35, 65	0
All	All	2591/2662 (97%)	-0.05	55 (2%) 63 59	11, 19, 40, 81	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1287	ASP	8.4
1	A	1323	ASN	7.1
1	A	1289	ALA	5.2
1	A	1317	VAL	4.6
1	A	61	LEU	4.4
1	A	1111	GLY	4.4
1	A	1324	CYS	4.3
1	B	61	LEU	4.2
1	B	218	LEU	4.1
1	B	1110	THR	3.9
1	B	220	ASP	3.8
1	B	1288	ASN	3.7
1	A	192	PRO	3.6
1	B	1287	ASP	3.6
1	A	529	ALA	3.5
1	B	1111	GLY	3.4
1	A	1322	GLU	3.4
1	A	1288	ASN	3.3
1	B	221	THR	3.3
1	B	565	LYS	3.2
1	B	1289	ALA	3.2
1	A	473	LEU	3.1
1	A	1110	THR	3.1
1	A	3	ALA	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	1106	LYS	3.1
1	A	220	ASP	3.0
1	A	534	MET	2.9
1	A	218	LEU	2.8
1	A	221	THR	2.8
1	A	386	HIS	2.8
1	A	530	ASP	2.7
1	A	566	ASP	2.6
1	A	1325	LYS	2.6
1	A	744	LEU	2.5
1	A	387	THR	2.5
1	B	3	ALA	2.5
1	A	531	LEU	2.5
1	A	537	LYS	2.4
1	A	222	PRO	2.4
1	A	474	SER	2.4
1	A	448	VAL	2.4
1	A	565	LYS	2.4
1	B	222	PRO	2.4
1	A	1229	ILE	2.4
1	B	1320	VAL	2.4
1	B	566	ASP	2.3
1	A	383	ARG	2.3
1	A	742	PHE	2.2
1	B	473	LEU	2.2
1	A	724	SER	2.2
1	A	1109	PRO	2.1
1	B	535	CYS	2.1
1	A	686	LYS	2.0
1	A	166	GLY	2.0
1	A	470	PRO	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 monosaccharides in this entry.

6.4 Ligands

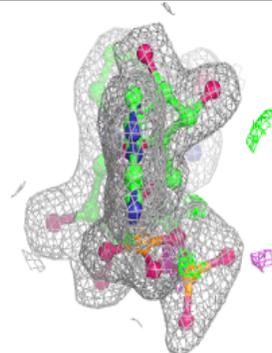
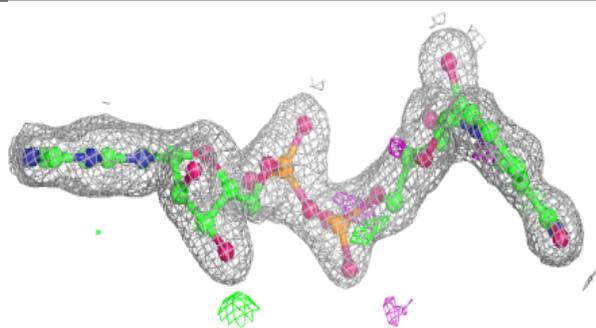
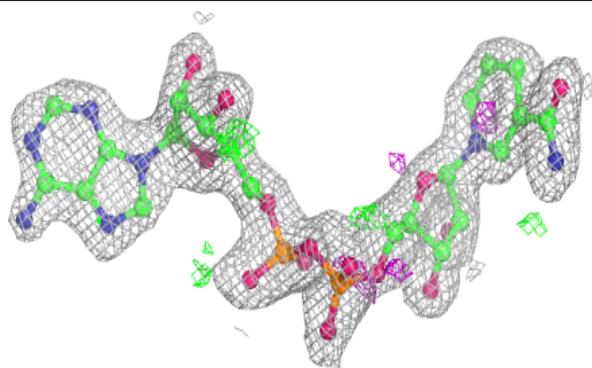
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	URC	A	1501	12/12	0.87	0.16	20,25,29,29	0
2	URC	B	3003	12/12	0.91	0.13	18,22,26,26	0
6	NAI	A	1506	44/44	0.92	0.10	23,32,39,43	0
6	NAI	B	3006	44/44	0.93	0.10	20,23,30,35	0
5	FAD	A	1505	53/53	0.96	0.09	17,20,24,25	0
3	BCT	A	1502	4/4	0.98	0.09	18,21,23,25	0
5	FAD	B	3005	53/53	0.98	0.08	13,14,16,17	0
4	FES	A	1504	4/4	0.99	0.05	16,17,17,18	0
4	FES	B	3001	4/4	0.99	0.05	15,15,17,18	0
3	BCT	B	3004	4/4	0.99	0.12	14,15,17,19	0
4	FES	A	1503	4/4	1.00	0.09	13,14,14,14	0
4	FES	B	3002	4/4	1.00	0.08	11,11,11,12	0

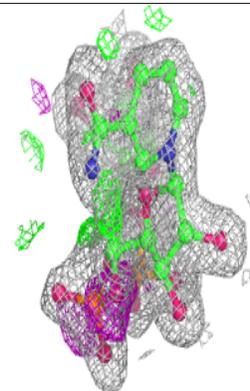
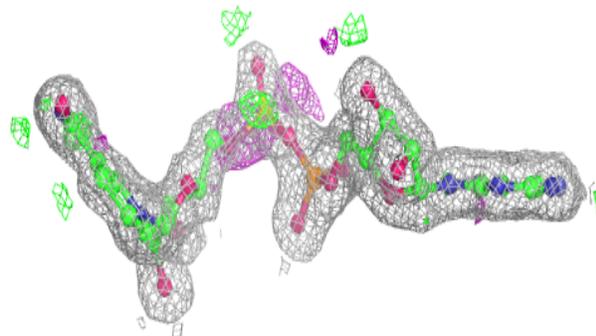
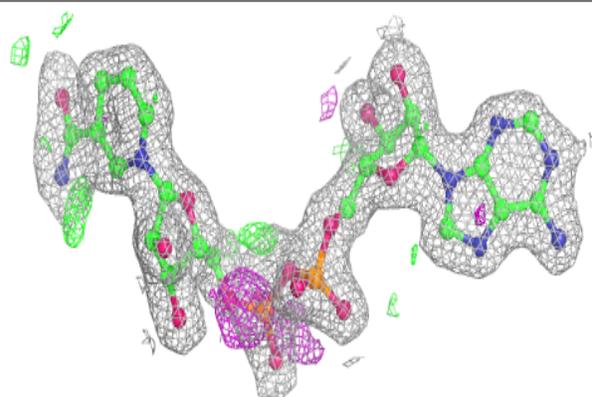
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAI A 1506:

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

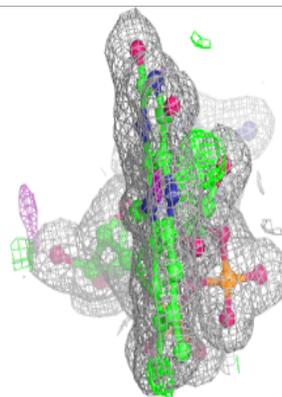
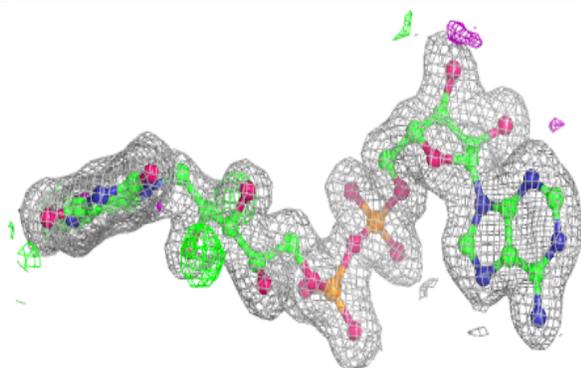
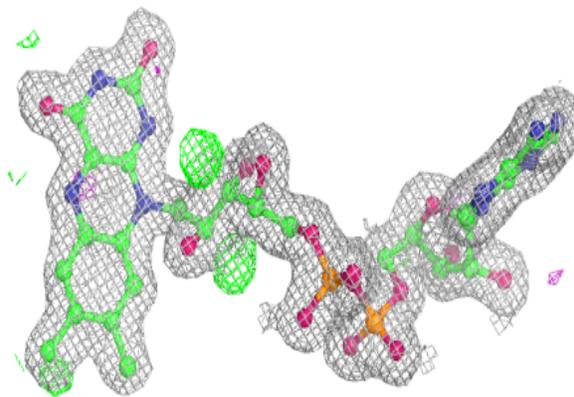
**Electron density around NAI B 3006:**

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

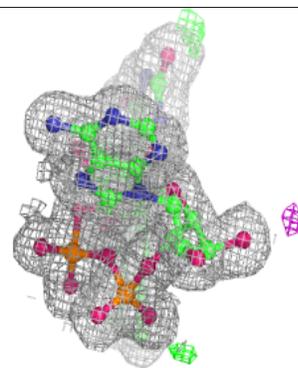
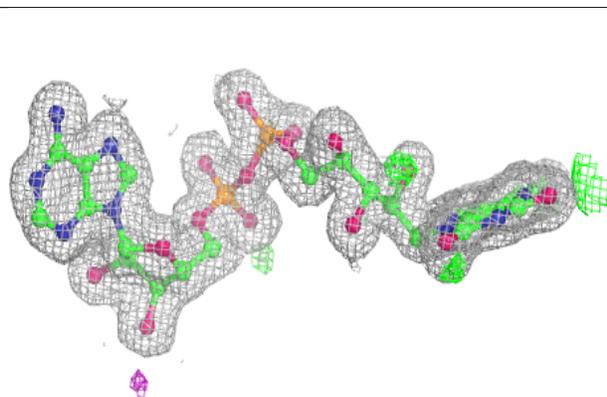
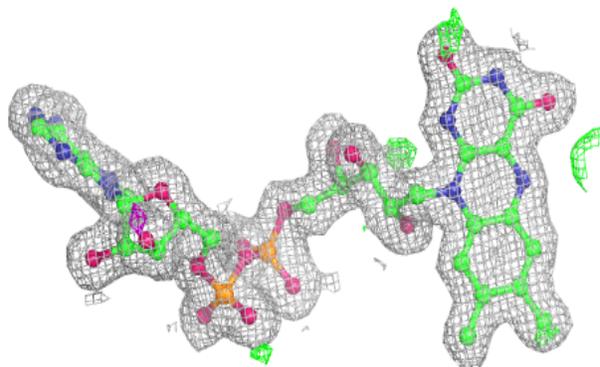


Electron density around FAD A 1505:

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

**Electron density around FAD B 3005:**

$2mF_o-DF_c$ (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.