



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:29 PM UTC

PDB ID : 5VWT / pdb_00005vwt
Title : Crystal structure of oxidized Aspergillus fumigatus UDP-galactopyranose mutase complexed with NADPH
Authors : Tanner, J.J.
Deposited on : 2017-05-22
Resolution : 2.75 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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

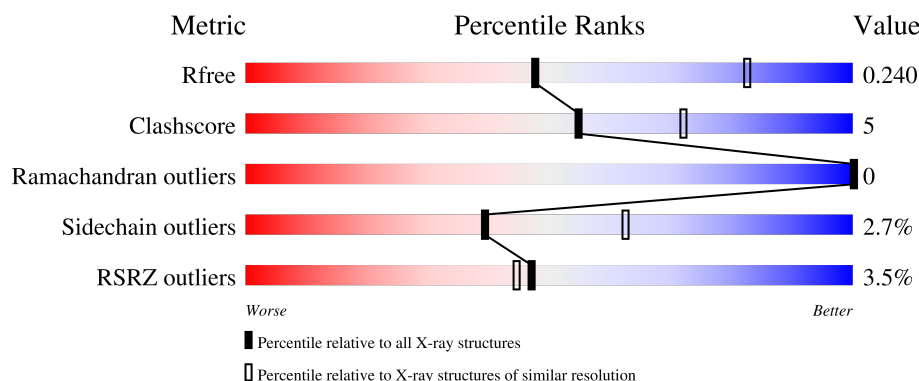
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	513	2% 84% 12% .
1	B	513	3% 87% 9% ..
1	C	513	5% 87% 10% .
1	D	513	4% 85% 11% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15472 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 UDP-galactopyranose mutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	495	Total	C	N	O	S	0	0	0
			3756	2396	632	707	21			
1	B	492	Total	C	N	O	S	0	0	0
			3748	2384	633	711	20			
1	C	501	Total	C	N	O	S	0	0	0
			3774	2400	639	716	19			
1	D	499	Total	C	N	O	S	0	0	0
			3740	2380	629	713	18			

There are 24 discrepancies between the modelled and reference sequences:

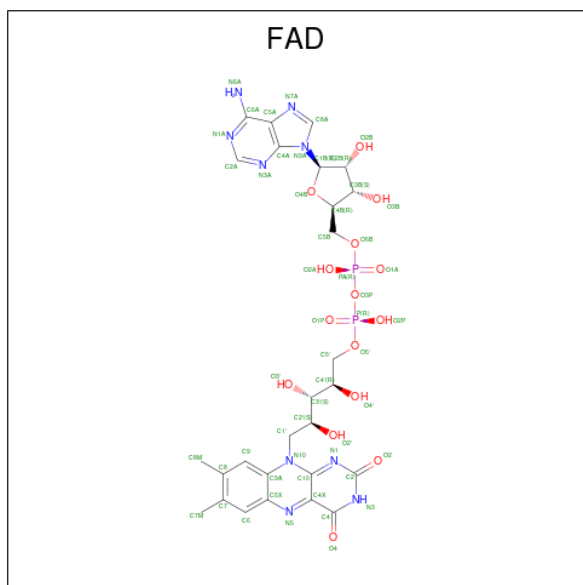
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	ALA	-	expression tag	UNP Q4W1X2
A	-1	ILE	-	expression tag	UNP Q4W1X2
A	0	ALA	-	expression tag	UNP Q4W1X2
A	344	ALA	LYS	engineered mutation	UNP Q4W1X2
A	345	ALA	LYS	engineered mutation	UNP Q4W1X2
A	429	THR	ALA	conflict	UNP Q4W1X2
B	-2	ALA	-	expression tag	UNP Q4W1X2
B	-1	ILE	-	expression tag	UNP Q4W1X2
B	0	ALA	-	expression tag	UNP Q4W1X2
B	344	ALA	LYS	engineered mutation	UNP Q4W1X2
B	345	ALA	LYS	engineered mutation	UNP Q4W1X2
B	429	THR	ALA	conflict	UNP Q4W1X2
C	-2	ALA	-	expression tag	UNP Q4W1X2
C	-1	ILE	-	expression tag	UNP Q4W1X2
C	0	ALA	-	expression tag	UNP Q4W1X2
C	344	ALA	LYS	engineered mutation	UNP Q4W1X2
C	345	ALA	LYS	engineered mutation	UNP Q4W1X2
C	429	THR	ALA	conflict	UNP Q4W1X2
D	-2	ALA	-	expression tag	UNP Q4W1X2
D	-1	ILE	-	expression tag	UNP Q4W1X2
D	0	ALA	-	expression tag	UNP Q4W1X2

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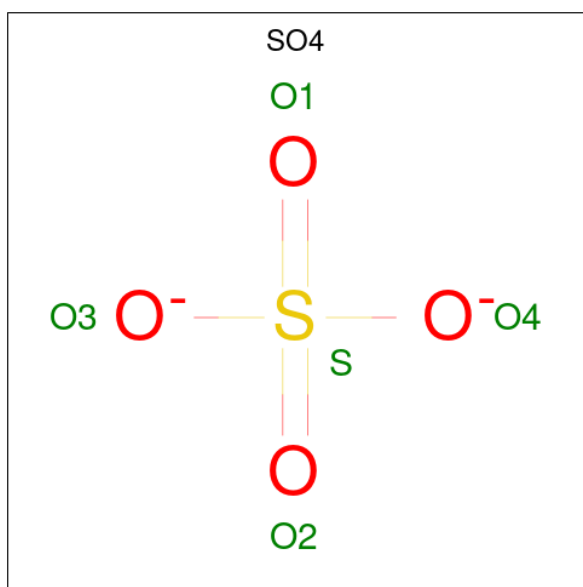
Chain	Residue	Modelled	Actual	Comment	Reference
D	344	ALA	LYS	engineered mutation	UNP Q4W1X2
D	345	ALA	LYS	engineered mutation	UNP Q4W1X2
D	429	THR	ALA	conflict	UNP Q4W1X2

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



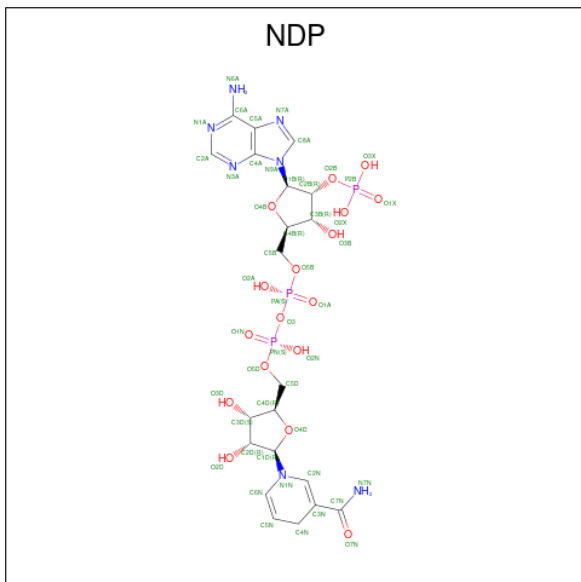
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



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

- Molecule 4 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: $\text{C}_{21}\text{H}_{30}\text{N}_7\text{O}_{17}\text{P}_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 48	C 21	N 7	O 17	P 3	0	0
4	B	1	Total 48	C 21	N 7	O 17	P 3	0	0
4	C	1	Total 31	C 10	N 5	O 13	P 3	0	0
4	D	1	Total 31	C 10	N 5	O 13	P 3	0	0

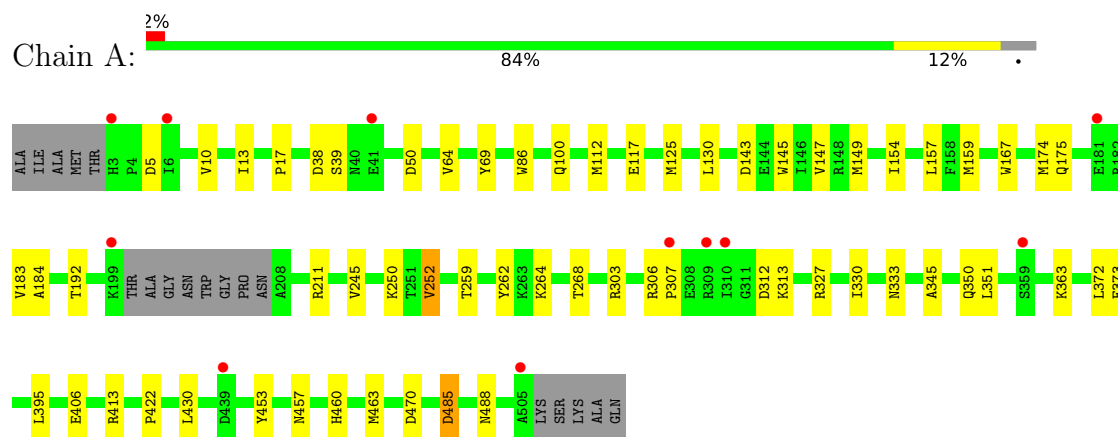
- Molecule 5 is water.

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

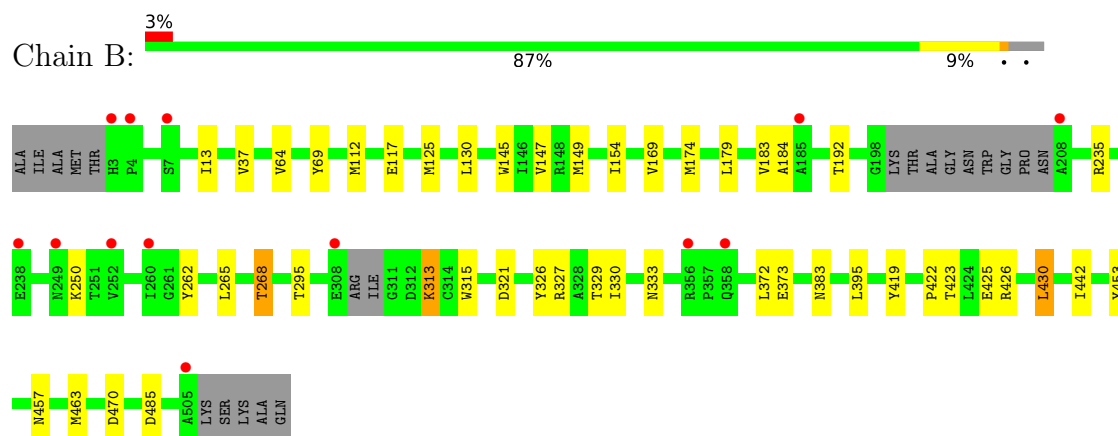
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

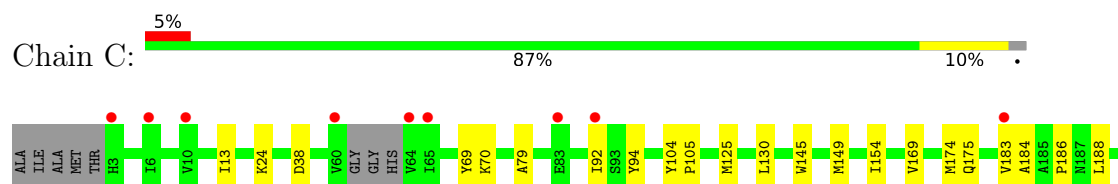
• Molecule 1: UDP-galactopyranose mutase

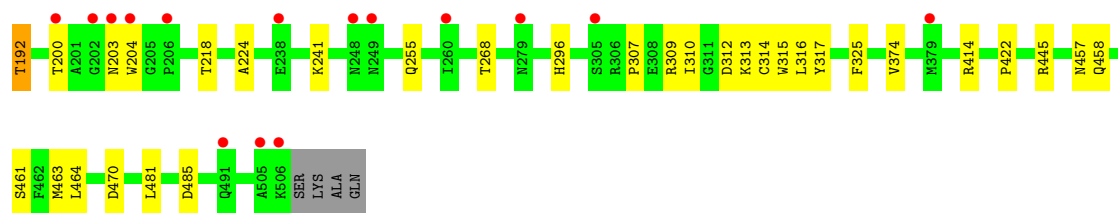


• Molecule 1: UDP-galactopyranose mutase

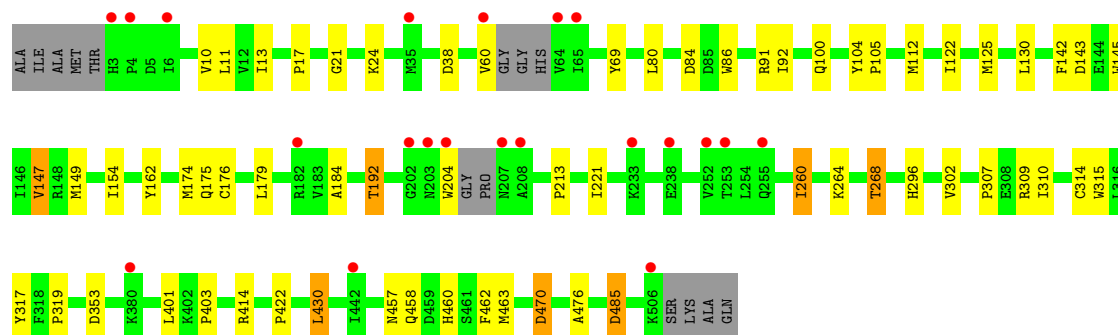
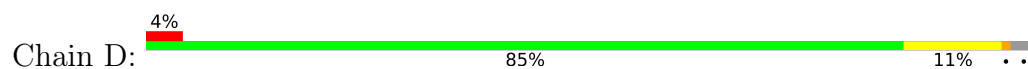


• Molecule 1: UDP-galactopyranose mutase





● Molecule 1: UDP-galactopyranose mutase



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	218.29Å 218.29Å 319.22Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	162.66 – 2.75 162.66 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.8 (162.66-2.75) 99.7 (162.66-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 2.73Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.202 , 0.239 0.203 , 0.240	Depositor DCC
R_{free} test set	5839 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	46.9	Xtriage
Anisotropy	0.460	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 57.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	15472	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage'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: FAD, SO4, NDP

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	0.40	0/3850	0.59	0/5255
1	B	0.38	0/3840	0.58	1/5240 (0.0%)
1	C	0.37	0/3869	0.56	0/5287
1	D	0.37	0/3834	0.56	0/5243
All	All	0.38	0/15393	0.57	1/21025 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	321	ASP	CB-CA-C	-5.04	99.88	110.17

There are no chirality outliers.

There are no planarity outliers.

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	3756	0	3538	35	0
1	B	3748	0	3543	27	0
1	C	3774	0	3525	32	0
1	D	3740	0	3444	44	0
2	A	53	0	31	7	0
2	B	53	0	31	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	53	0	31	6	0
2	D	53	0	31	7	0
3	A	15	0	0	0	0
3	B	20	0	0	1	0
3	C	20	0	0	0	0
3	D	15	0	0	1	0
4	A	48	0	26	2	0
4	B	48	0	26	1	0
4	C	31	0	11	0	0
4	D	31	0	11	0	0
5	A	7	0	0	0	0
5	B	4	0	0	0	0
5	C	2	0	0	0	0
5	D	1	0	0	1	0
All	All	15472	0	14248	142	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 5.

All (142) 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:D:17:PRO:HD2	2:D:601:FAD:H5'2	1.57	0.86
1:A:17:PRO:HD2	2:A:601:FAD:H5'2	1.58	0.84
1:A:372:LEU:HD21	1:A:395:LEU:HD21	1.65	0.79
1:B:13:ILE:O	1:B:268:THR:HB	1.88	0.73
1:A:268:THR:HG22	2:A:601:FAD:H51A	1.71	0.71
1:A:13:ILE:O	1:A:268:THR:HB	1.92	0.69
1:A:174:MET:HE3	1:A:422:PRO:HD2	1.78	0.65
1:A:245:VAL:HG22	1:A:252:VAL:HG22	1.78	0.65
1:D:307:PRO:HG2	1:D:310:ILE:HD13	1.79	0.64
1:D:125:MET:HE3	1:D:154:ILE:HD12	1.80	0.64
1:C:307:PRO:HB2	1:C:309:ARG:HG2	1.78	0.64
1:B:313:LYS:H	1:B:313:LYS:HD2	1.63	0.63
1:B:327:ARG:NE	1:B:373:GLU:OE1	2.22	0.62
1:B:313:LYS:HG3	1:B:333:ASN:HB3	1.80	0.61
1:A:149:MET:SD	1:A:184:ALA:HA	2.41	0.61
1:D:143:ASP:O	1:D:147:VAL:HG23	2.00	0.60
1:B:149:MET:SD	1:B:184:ALA:HA	2.41	0.60
1:C:315:TRP:CD1	1:C:317:TYR:HH	2.19	0.60
1:B:145:TRP:CZ2	1:B:149:MET:HE2	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:69:TYR:CD2	1:D:463:MET:HG3	2.36	0.59
1:B:372:LEU:HD21	1:B:395:LEU:HD21	1.83	0.59
1:D:309:ARG:NH1	3:D:603:SO4:O3	2.29	0.59
1:D:174:MET:HE3	1:D:422:PRO:HD2	1.82	0.59
1:B:174:MET:HE3	1:B:422:PRO:HD2	1.85	0.58
1:C:13:ILE:O	1:C:268:THR:HB	2.04	0.58
1:D:470:ASP:HB3	1:D:476:ALA:HB3	1.87	0.56
1:D:13:ILE:O	1:D:268:THR:HB	2.06	0.56
1:A:192:THR:HG21	1:D:122:ILE:HG21	1.88	0.55
1:C:38:ASP:OD1	2:C:601:FAD:H1B	2.06	0.55
1:C:125:MET:HE3	1:C:154:ILE:HD12	1.87	0.55
1:B:383:ASN:ND2	3:B:605:SO4:O4	2.34	0.54
1:A:125:MET:HE3	1:A:154:ILE:HD12	1.90	0.54
1:B:130:LEU:HD23	1:C:130:LEU:HD23	1.91	0.54
1:D:457:ASN:HB2	2:D:601:FAD:O2	2.09	0.53
1:C:268:THR:HG22	2:C:601:FAD:H52A	1.91	0.53
1:D:92:ILE:O	1:D:314:CYS:HB2	2.09	0.52
1:B:265:LEU:HD23	1:B:442:ILE:HG12	1.92	0.52
1:C:188:LEU:O	1:C:192:THR:HB	2.10	0.52
1:D:69:TYR:CG	1:D:463:MET:HG3	2.45	0.52
1:A:268:THR:CG2	2:A:601:FAD:H51A	2.40	0.52
1:B:69:TYR:CG	1:B:463:MET:HG3	2.45	0.51
1:A:306:ARG:HB2	1:A:307:PRO:HD2	1.92	0.51
1:D:315:TRP:CD1	1:D:317:TYR:HH	2.29	0.51
1:B:457:ASN:HB2	2:B:601:FAD:O2	2.12	0.50
1:C:325:PHE:HA	1:C:374:VAL:HG22	1.94	0.50
1:D:80:LEU:HD12	1:D:86:TRP:CZ2	2.47	0.50
1:A:69:TYR:CD2	1:A:463:MET:HG3	2.47	0.50
1:B:453:TYR:CE1	4:B:603:NDP:H42N	2.47	0.50
1:B:313:LYS:HD2	1:B:313:LYS:N	2.26	0.49
1:A:457:ASN:HB3	4:A:604:NDP:H41N	1.94	0.49
1:D:100:GLN:HB3	1:D:112:MET:HE2	1.95	0.49
1:B:125:MET:HE3	1:B:154:ILE:HD12	1.94	0.49
1:D:92:ILE:HA	1:D:204:TRP:CE3	2.48	0.49
1:A:457:ASN:HB2	2:A:601:FAD:O2	2.13	0.49
1:C:69:TYR:CG	1:C:463:MET:HG3	2.48	0.48
1:A:5:ASP:HB2	1:A:259:THR:O	2.13	0.48
1:A:69:TYR:CG	1:A:463:MET:HG3	2.47	0.48
1:C:149:MET:SD	1:C:184:ALA:HA	2.54	0.48
1:C:174:MET:HE3	1:C:422:PRO:HD2	1.96	0.48
1:C:445:ARG:NE	1:C:481:LEU:HD22	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:458:GLN:HG3	2:C:601:FAD:O2	2.14	0.47
1:D:91:ARG:O	1:D:204:TRP:HB3	2.15	0.47
1:C:92:ILE:O	1:C:314:CYS:HB2	2.15	0.47
1:A:143:ASP:O	1:A:147:VAL:HG23	2.15	0.47
1:C:461:SER:HA	1:C:464:LEU:HD12	1.97	0.46
1:A:457:ASN:HB2	2:A:601:FAD:C2	2.45	0.46
1:A:485:ASP:OD1	1:A:485:ASP:N	2.42	0.46
1:A:250:LYS:HG2	1:A:262:TYR:CE1	2.51	0.46
1:B:145:TRP:CD2	1:B:179:LEU:HD13	2.50	0.46
1:D:192:THR:HG23	5:D:701:HOH:O	2.14	0.46
2:B:601:FAD:H2'	2:B:601:FAD:N1	2.31	0.46
1:A:86:TRP:CD2	1:A:211:ARG:HD2	2.50	0.46
1:D:69:TYR:CE2	1:D:463:MET:HG3	2.51	0.46
1:D:84:ASP:OD1	1:D:84:ASP:N	2.45	0.46
1:D:213:PRO:HG3	1:D:221:ILE:HG13	1.97	0.46
1:D:11:LEU:HD13	1:D:260:ILE:HG21	1.98	0.46
1:A:10:VAL:HA	1:A:264:LYS:O	2.17	0.45
1:A:38:ASP:OD1	1:A:39:SER:N	2.50	0.45
1:B:37:VAL:HG12	1:B:235:ARG:HB3	1.96	0.45
1:B:315:TRP:CE3	1:B:329:THR:HB	2.52	0.45
1:C:457:ASN:HB2	2:C:601:FAD:O2	2.17	0.45
1:B:295:THR:HG21	2:B:601:FAD:C7M	2.46	0.45
1:B:426:ARG:O	1:B:430:LEU:HB2	2.17	0.45
1:D:296:HIS:CE1	1:D:414:ARG:HD2	2.52	0.45
1:C:104:TYR:CG	1:C:105:PRO:HA	2.52	0.45
1:C:458:GLN:HA	2:C:601:FAD:O3'	2.17	0.45
1:C:145:TRP:CZ2	1:C:149:MET:HE2	2.52	0.44
1:D:457:ASN:HB2	2:D:601:FAD:C2	2.48	0.44
1:A:460:HIS:NE2	1:A:488:ASN:OD1	2.46	0.44
1:C:310:ILE:O	1:C:313:LYS:HG3	2.18	0.44
1:B:69:TYR:CD1	1:B:463:MET:HG3	2.52	0.44
1:D:142:PHE:HA	1:D:176:CYS:HB3	1.99	0.44
1:D:149:MET:SD	1:D:184:ALA:HA	2.57	0.44
1:A:117:GLU:OE2	1:A:157:LEU:HD21	2.18	0.44
1:C:92:ILE:HG22	1:C:94:TYR:HE1	1.82	0.44
2:D:601:FAD:H9	2:D:601:FAD:H1'1	1.78	0.44
2:B:601:FAD:H9	2:B:601:FAD:H1'1	1.63	0.43
1:D:38:ASP:OD1	2:D:601:FAD:H1B	2.18	0.43
1:C:92:ILE:HD12	1:C:312:ASP:O	2.18	0.43
1:B:423:THR:HB	1:B:425:GLU:OE1	2.17	0.43
1:C:296:HIS:CE1	1:C:414:ARG:HD2	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:601:FAD:H9	2:C:601:FAD:H1'1	1.78	0.43
1:D:10:VAL:HA	1:D:264:LYS:O	2.19	0.43
1:B:250:LYS:HB3	1:B:262:TYR:O	2.19	0.43
1:D:457:ASN:OD1	1:D:460:HIS:ND1	2.51	0.43
1:D:458:GLN:HG3	2:D:601:FAD:O2	2.19	0.43
1:D:353:ASP:HB3	1:D:403:PRO:HB3	2.01	0.43
1:A:100:GLN:HB3	1:A:112:MET:HE2	2.01	0.43
1:B:419:TYR:HE1	2:B:601:FAD:HM72	1.84	0.43
1:C:79:ALA:HB1	1:C:224:ALA:CB	2.48	0.42
2:A:601:FAD:O2'	2:A:601:FAD:H9	2.19	0.42
4:A:604:NDP:H71N	4:A:604:NDP:H2N	1.55	0.42
1:A:350:GLN:HG3	1:A:351:LEU:O	2.19	0.42
1:C:69:TYR:CD1	1:C:463:MET:HG3	2.54	0.42
1:A:327:ARG:NE	1:A:373:GLU:OE1	2.40	0.42
1:D:145:TRP:CD2	1:D:179:LEU:HD13	2.55	0.42
1:D:430:LEU:HD12	1:D:430:LEU:HA	1.88	0.42
1:C:313:LYS:HD2	1:C:316:LEU:HD21	2.02	0.42
1:D:307:PRO:HB2	1:D:309:ARG:HG2	2.02	0.42
1:A:50:ASP:OD2	1:A:413:ARG:NH1	2.47	0.41
1:B:112:MET:HB3	1:B:112:MET:HE2	1.78	0.41
1:D:162:TYR:HB2	1:D:319:PRO:HB3	2.01	0.41
1:D:24:LYS:HD2	1:D:24:LYS:HA	1.88	0.41
1:D:104:TYR:CG	1:D:105:PRO:HA	2.55	0.41
1:C:24:LYS:HD2	1:C:24:LYS:HA	1.90	0.41
1:A:130:LEU:HD23	1:D:130:LEU:HD23	2.03	0.41
1:C:241:LYS:HB3	1:C:255:GLN:HB2	2.02	0.41
1:A:145:TRP:CZ2	1:A:149:MET:HE2	2.56	0.40
1:A:313:LYS:HG3	1:A:333:ASN:HB3	2.03	0.40
1:A:345:ALA:HA	1:A:363:LYS:O	2.21	0.40
1:B:326:TYR:CZ	1:B:373:GLU:HB3	2.56	0.40
1:D:302:VAL:HG21	1:D:401:LEU:HD21	2.02	0.40
1:A:303:ARG:NE	1:A:406:GLU:OE1	2.50	0.40
2:A:601:FAD:HM71	2:A:601:FAD:HM83	1.90	0.40
1:D:17:PRO:CD	2:D:601:FAD:H5'2	2.40	0.40
1:D:21:GLY:HA2	1:D:462:PHE:CE1	2.56	0.40
1:D:485:ASP:OD1	1:D:485:ASP:N	2.43	0.40
1:C:92:ILE:HA	1:C:204:TRP:CE3	2.56	0.40
1:C:105:PRO:HD3	1:C:203:ASN:HB3	2.03	0.40
1:C:149:MET:O	1:C:186:PRO:HB3	2.20	0.40
1:A:167:TRP:HH2	1:A:453:TYR:CG	2.39	0.40
1:D:142:PHE:CE1	1:D:179:LEU:HD11	2.56	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	491/513 (96%)	479 (98%)	12 (2%)	0	100	100
1	B	486/513 (95%)	470 (97%)	16 (3%)	0	100	100
1	C	497/513 (97%)	482 (97%)	15 (3%)	0	100	100
1	D	493/513 (96%)	478 (97%)	15 (3%)	0	100	100
All	All	1967/2052 (96%)	1909 (97%)	58 (3%)	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	378/432 (88%)	368 (97%)	10 (3%)	40	63
1	B	383/432 (89%)	371 (97%)	12 (3%)	35	59
1	C	378/432 (88%)	369 (98%)	9 (2%)	43	65
1	D	368/432 (85%)	359 (98%)	9 (2%)	43	65
All	All	1507/1728 (87%)	1467 (97%)	40 (3%)	39	62

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	VAL
1	A	159	MET
1	A	175	GLN
1	A	183	VAL
1	A	252	VAL
1	A	312	ASP
1	A	330	ILE
1	A	430	LEU
1	A	470	ASP
1	A	485	ASP
1	B	64	VAL
1	B	117	GLU
1	B	147	VAL
1	B	169	VAL
1	B	183	VAL
1	B	192	THR
1	B	268	THR
1	B	313	LYS
1	B	330	ILE
1	B	430	LEU
1	B	470	ASP
1	B	485	ASP
1	C	70	LYS
1	C	169	VAL
1	C	175	GLN
1	C	183	VAL
1	C	192	THR
1	C	200	THR
1	C	218	THR
1	C	470	ASP
1	C	485	ASP
1	D	60	VAL
1	D	147	VAL
1	D	175	GLN
1	D	192	THR
1	D	260	ILE
1	D	268	THR
1	D	430	LEU
1	D	470	ASP
1	D	485	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	GLN
1	A	246	ASN
1	B	27	ASN
1	B	175	GLN
1	B	279	ASN
1	C	98	GLN
1	D	40	ASN
1	D	98	GLN
1	D	393	GLN

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](#)

22 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	SO4	A	602	-	4,4,4	0.30	0	6,6,6	0.34	0
4	NDP	C	602	-	32,33,52	3.93	8 (25%)	50,52,80	1.63	12 (24%)
3	SO4	D	605	-	4,4,4	0.22	0	6,6,6	0.27	0
3	SO4	B	605	-	4,4,4	0.28	0	6,6,6	0.15	0
3	SO4	D	603	-	4,4,4	0.28	0	6,6,6	0.26	0
3	SO4	C	606	-	4,4,4	0.23	0	6,6,6	0.23	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FAD	B	601	-	58,58,58	0.55	1 (1%)	85,89,89	0.78	2 (2%)
3	SO4	D	602	-	4,4,4	0.29	0	6,6,6	0.45	0
3	SO4	B	604	-	4,4,4	0.22	0	6,6,6	0.34	0
3	SO4	A	605	-	4,4,4	0.34	0	6,6,6	0.44	0
2	FAD	A	601	-	58,58,58	0.42	0	85,89,89	0.81	3 (3%)
3	SO4	C	604	-	4,4,4	0.26	0	6,6,6	0.20	0
3	SO4	A	603	-	4,4,4	0.28	0	6,6,6	0.12	0
4	NDP	A	604	-	51,52,52	2.95	15 (29%)	71,80,80	1.46	14 (19%)
2	FAD	C	601	-	58,58,58	0.34	0	85,89,89	0.48	0
4	NDP	D	604	-	32,33,52	3.66	8 (25%)	50,52,80	1.74	14 (28%)
3	SO4	B	606	-	4,4,4	0.27	0	6,6,6	0.28	0
3	SO4	C	603	-	4,4,4	0.27	0	6,6,6	0.38	0
2	FAD	D	601	-	58,58,58	0.41	0	85,89,89	0.61	1 (1%)
3	SO4	C	605	-	4,4,4	0.25	0	6,6,6	0.21	0
4	NDP	B	603	-	51,52,52	3.05	11 (21%)	71,80,80	1.39	13 (18%)
3	SO4	B	602	-	4,4,4	0.28	0	6,6,6	0.39	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
2	FAD	C	601	-	-	12/34/50/50	0/6/6/6
4	NDP	D	604	-	-	3/21/37/77	0/3/3/5
4	NDP	B	603	-	-	2/34/77/77	0/5/5/5
4	NDP	C	602	-	-	7/21/37/77	0/3/3/5
2	FAD	A	601	-	-	15/34/50/50	0/6/6/6
2	FAD	D	601	-	-	14/34/50/50	0/6/6/6
2	FAD	B	601	-	-	13/34/50/50	0/6/6/6
4	NDP	A	604	-	-	5/34/77/77	0/5/5/5

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	602	NDP	P2B-O2B	17.23	1.89	1.59
4	B	603	NDP	P2B-O2B	16.24	1.88	1.59
4	D	604	NDP	P2B-O2B	15.30	1.86	1.59
4	A	604	NDP	P2B-O2B	13.66	1.83	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	604	NDP	PA-O3	11.16	1.71	1.59
4	B	603	NDP	PA-O3	9.53	1.69	1.59
4	C	602	NDP	PA-O3	9.25	1.69	1.59
4	D	604	NDP	PN-O1N	8.33	1.76	1.50
4	D	604	NDP	PA-O3	7.60	1.67	1.59
4	C	602	NDP	PN-O5D	7.55	1.82	1.54
4	A	604	NDP	PN-O5D	5.48	1.80	1.59
4	B	603	NDP	PN-O5D	4.52	1.77	1.59
4	B	603	NDP	PN-O3	3.64	1.63	1.59
4	A	604	NDP	PN-O3	3.29	1.63	1.59
4	A	604	NDP	O2B-C2B	-3.17	1.33	1.44
4	D	604	NDP	C8A-N9A	3.11	1.42	1.37
4	A	604	NDP	C3D-C4D	2.99	1.60	1.53
4	B	603	NDP	C2A-N1A	2.92	1.39	1.33
4	B	603	NDP	C5A-C4A	2.90	1.44	1.39
2	B	601	FAD	P-O3P	-2.88	1.56	1.59
4	D	604	NDP	O2B-C2B	-2.85	1.34	1.44
4	D	604	NDP	C5A-C4A	2.83	1.44	1.39
4	A	604	NDP	C8A-N9A	2.75	1.42	1.37
4	D	604	NDP	C8A-N7A	2.72	1.36	1.31
4	B	603	NDP	O2B-C2B	-2.70	1.34	1.44
4	A	604	NDP	C5A-C4A	2.67	1.43	1.39
4	C	602	NDP	O2B-C2B	-2.64	1.35	1.44
4	C	602	NDP	C8A-N9A	2.56	1.41	1.37
4	B	603	NDP	C4A-N3A	2.52	1.39	1.34
4	D	604	NDP	C2A-N1A	2.51	1.38	1.33
4	B	603	NDP	C6N-N1N	2.48	1.43	1.37
4	B	603	NDP	C8A-N9A	2.45	1.41	1.37
4	A	604	NDP	C4A-N3A	2.40	1.38	1.34
4	A	604	NDP	C2A-N1A	2.33	1.38	1.33
4	C	602	NDP	C2A-N1A	2.33	1.38	1.33
4	C	602	NDP	C5A-C4A	2.22	1.43	1.39
4	A	604	NDP	C2B-C1B	2.18	1.58	1.53
4	A	604	NDP	C8A-N7A	2.14	1.35	1.31
4	A	604	NDP	C7N-C3N	-2.12	1.44	1.48
4	C	602	NDP	C8A-N7A	2.10	1.35	1.31
4	A	604	NDP	C7N-N7N	2.07	1.39	1.33
4	A	604	NDP	PA-O5B	2.04	1.67	1.59
4	B	603	NDP	C3D-C4D	2.03	1.58	1.53

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	FAD	O3'-C3'-C2'	3.86	117.69	108.93
4	D	604	NDP	O5D-PN-O2N	3.84	122.21	107.80
4	A	604	NDP	O2B-P2B-O1X	-3.84	95.65	109.33
4	D	604	NDP	O2N-PN-O3	3.66	116.90	104.64
4	A	604	NDP	P2B-O2B-C2B	-3.45	114.21	123.43
4	C	602	NDP	P2B-O2B-C2B	-3.44	114.24	123.43
4	D	604	NDP	P2B-O2B-C2B	-3.37	114.44	123.43
4	D	604	NDP	PA-O5B-C5B	-3.10	103.61	121.35
4	D	604	NDP	O5D-PN-O1N	-3.09	98.79	110.83
4	C	602	NDP	PA-O5B-C5B	-3.03	103.96	121.35
4	C	602	NDP	O3X-P2B-O2B	-2.96	94.30	105.85
4	C	602	NDP	C3B-C2B-C1B	-2.96	97.14	102.81
4	D	604	NDP	C3B-C2B-C1B	-2.93	97.20	102.81
4	B	603	NDP	P2B-O2B-C2B	-2.87	115.77	123.43
4	C	602	NDP	O2B-P2B-O1X	-2.81	99.31	109.33
4	C	602	NDP	O5D-PN-O1N	-2.81	99.89	110.83
4	A	604	NDP	PA-O5B-C5B	-2.77	105.48	121.35
2	D	601	FAD	C4'-C3'-C2'	2.76	118.16	113.57
4	C	602	NDP	O2N-PN-O1N	2.69	121.30	110.83
4	C	602	NDP	O5D-PN-O3	2.63	113.45	104.64
4	D	604	NDP	O3X-P2B-O2X	2.61	117.58	107.80
4	B	603	NDP	O3X-P2B-O2B	-2.59	95.76	105.85
4	A	604	NDP	O3X-P2B-O2X	2.57	117.42	107.80
4	C	602	NDP	O3X-P2B-O2X	2.55	117.38	107.80
4	B	603	NDP	O3X-P2B-O2X	2.55	117.36	107.80
4	B	603	NDP	PA-O5B-C5B	-2.51	106.94	121.35
4	A	604	NDP	O2A-PA-O3	-2.46	100.64	107.27
4	D	604	NDP	C2A-N1A-C6A	-2.43	114.73	118.73
4	B	603	NDP	N3A-C4A-N9A	2.43	131.30	127.17
4	B	603	NDP	O2A-PA-O3	-2.42	100.74	107.27
4	D	604	NDP	O2A-PA-O3	-2.39	100.82	107.27
2	A	601	FAD	C4'-C3'-C2'	2.34	117.47	113.57
4	B	603	NDP	O4D-C1D-N1N	2.34	112.55	108.08
4	D	604	NDP	O2X-P2B-O2B	-2.33	96.78	105.85
4	B	603	NDP	O2B-P2B-O1X	-2.30	101.12	109.33
4	B	603	NDP	O7N-C7N-C3N	2.25	125.13	120.90
4	A	604	NDP	O4B-C4B-C3B	2.24	109.61	105.15
4	A	604	NDP	N3A-C4A-N9A	2.24	130.97	127.17
4	C	602	NDP	C5B-C4B-C3B	-2.22	107.23	115.21
4	D	604	NDP	O2N-PN-O1N	-2.21	102.23	110.83
4	A	604	NDP	C2A-N1A-C6A	-2.19	115.13	118.73
2	A	601	FAD	O3'-C3'-C4'	-2.18	103.98	108.93
4	A	604	NDP	O4D-C1D-N1N	2.16	112.19	108.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	604	NDP	C5A-C4A-N3A	-2.15	123.75	126.72
4	D	604	NDP	O4B-C4B-C3B	2.12	109.36	105.15
4	B	603	NDP	O2N-PN-O1N	2.11	122.27	112.44
4	B	603	NDP	C5A-C4A-N3A	-2.11	123.81	126.72
4	B	603	NDP	C2A-N1A-C6A	-2.11	115.27	118.73
4	C	602	NDP	O4B-C4B-C3B	2.10	109.32	105.15
4	A	604	NDP	O3-PA-O1A	-2.09	104.41	110.70
2	B	601	FAD	O3P-P-O1P	-2.07	104.47	110.70
4	A	604	NDP	C3N-C2N-N1N	-2.06	120.17	123.20
4	D	604	NDP	C5B-C4B-C3B	-2.05	107.83	115.21
4	A	604	NDP	C2D-C1D-N1N	-2.04	108.28	113.31
4	B	603	NDP	C3B-C2B-C1B	-2.04	98.89	102.81
2	A	601	FAD	O4'-C4'-C5'	2.03	114.47	109.99
4	A	604	NDP	C1D-N1N-C2N	-2.03	117.80	121.14
4	D	604	NDP	C2B-C1B-N9A	2.02	117.08	113.75
4	C	602	NDP	C2A-N1A-C6A	-2.02	115.42	118.73

There are no chirality outliers.

All (71) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	FAD	C2'-C1'-N10-C10
2	A	601	FAD	N10-C1'-C2'-O2'
2	A	601	FAD	C2'-C3'-C4'-O4'
2	A	601	FAD	C2'-C3'-C4'-C5'
2	A	601	FAD	O3'-C3'-C4'-O4'
2	A	601	FAD	O3'-C3'-C4'-C5'
2	A	601	FAD	C3'-C4'-C5'-O5'
2	A	601	FAD	O4'-C4'-C5'-O5'
2	A	601	FAD	C5'-O5'-P-O1P
2	A	601	FAD	C5'-O5'-P-O3P
2	B	601	FAD	C5B-O5B-PA-O1A
2	B	601	FAD	C5B-O5B-PA-O3P
2	B	601	FAD	P-O3P-PA-O5B
2	B	601	FAD	C1'-C2'-C3'-O3'
2	B	601	FAD	C1'-C2'-C3'-C4'
2	B	601	FAD	O2'-C2'-C3'-O3'
2	B	601	FAD	O2'-C2'-C3'-C4'
2	B	601	FAD	C3'-C4'-C5'-O5'
2	B	601	FAD	O4'-C4'-C5'-O5'
2	C	601	FAD	C1'-C2'-C3'-O3'
2	C	601	FAD	C1'-C2'-C3'-C4'

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Mol	Chain	Res	Type	Atoms
2	C	601	FAD	O2'-C2'-C3'-O3'
2	C	601	FAD	O2'-C2'-C3'-C4'
2	C	601	FAD	O4'-C4'-C5'-O5'
2	C	601	FAD	C5'-O5'-P-O1P
2	C	601	FAD	C5'-O5'-P-O2P
2	C	601	FAD	C5'-O5'-P-O3P
2	C	601	FAD	PA-O3P-P-O5'
2	D	601	FAD	C5B-O5B-PA-O2A
2	D	601	FAD	C5B-O5B-PA-O3P
2	D	601	FAD	P-O3P-PA-O5B
2	D	601	FAD	C1'-C2'-C3'-O3'
2	D	601	FAD	C1'-C2'-C3'-C4'
2	D	601	FAD	O2'-C2'-C3'-O3'
2	D	601	FAD	O2'-C2'-C3'-C4'
2	D	601	FAD	C3'-C4'-C5'-O5'
2	D	601	FAD	O4'-C4'-C5'-O5'
4	A	604	NDP	PA-O3-PN-O5D
4	C	602	NDP	C5B-O5B-PA-O1A
4	C	602	NDP	C5B-O5B-PA-O3
4	C	602	NDP	PN-O3-PA-O5B
4	C	602	NDP	O4B-C4B-C5B-O5B
4	C	602	NDP	C3B-C4B-C5B-O5B
2	D	601	FAD	O4B-C4B-C5B-O5B
4	A	604	NDP	C3D-C4D-C5D-O5D
2	D	601	FAD	C3B-C4B-C5B-O5B
4	A	604	NDP	O4D-C4D-C5D-O5D
4	A	604	NDP	O4D-C1D-N1N-C2N
4	B	603	NDP	PA-O3-PN-O5D
2	A	601	FAD	P-O3P-PA-O1A
4	B	603	NDP	O4D-C1D-N1N-C2N
2	A	601	FAD	N10-C1'-C2'-C3'
2	A	601	FAD	C5'-O5'-P-O2P
2	B	601	FAD	C5B-O5B-PA-O2A
2	B	601	FAD	C2'-C1'-N10-C10
4	C	602	NDP	C5B-O5B-PA-O2A
2	D	601	FAD	C4'-C5'-O5'-P
4	D	604	NDP	C2B-O2B-P2B-O2X
2	B	601	FAD	O4B-C4B-C5B-O5B
4	A	604	NDP	C2N-C3N-C7N-N7N
4	D	604	NDP	O4B-C4B-C5B-O5B
2	C	601	FAD	PA-O3P-P-O1P
2	D	601	FAD	O3'-C3'-C4'-C5'

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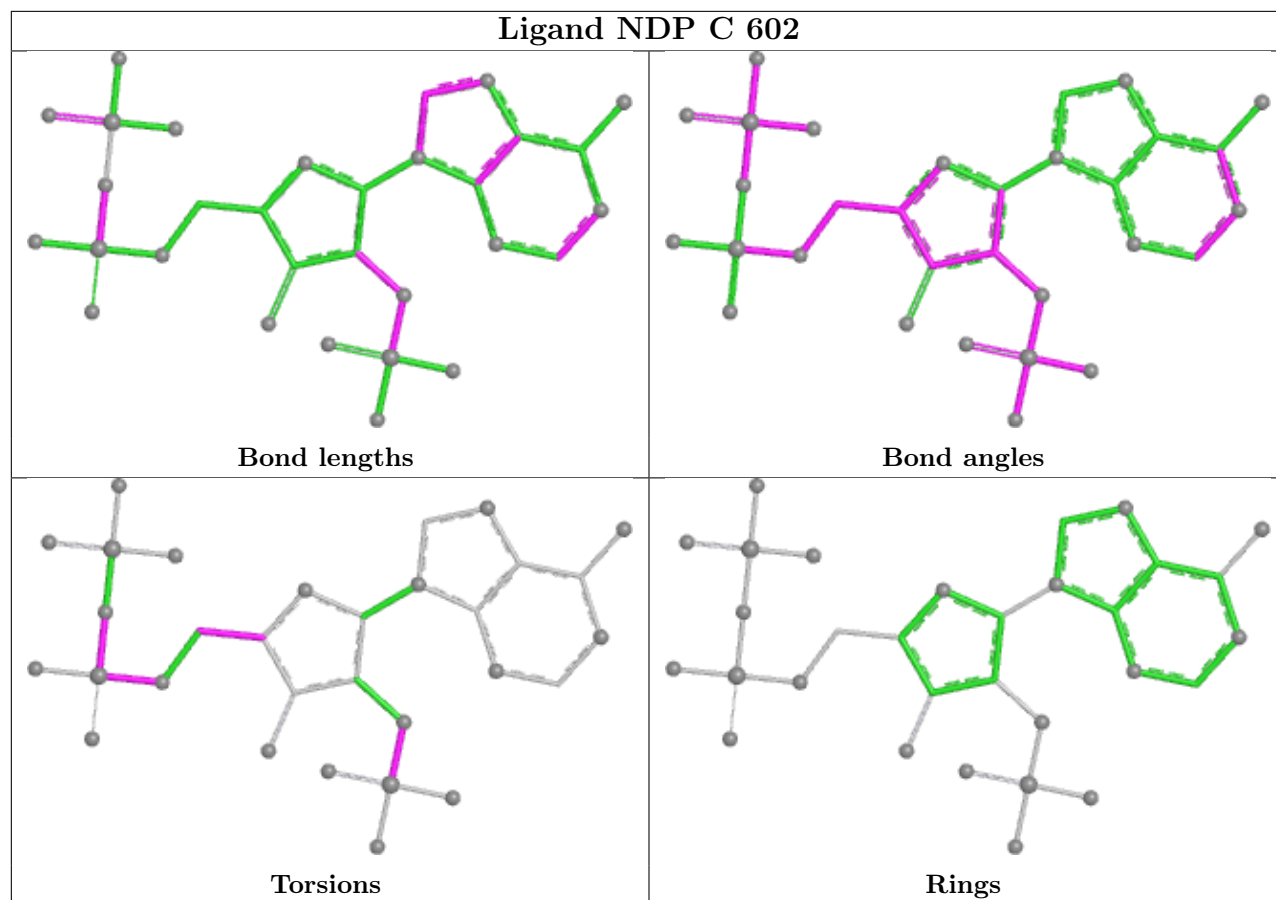
Mol	Chain	Res	Type	Atoms
2	A	601	FAD	O4B-C4B-C5B-O5B
2	C	601	FAD	C3'-C4'-C5'-O5'
2	C	601	FAD	PA-O3P-P-O2P
4	C	602	NDP	C2B-O2B-P2B-O2X
4	D	604	NDP	C2B-O2B-P2B-O1X
2	B	601	FAD	C2'-C3'-C4'-O4'
2	A	601	FAD	P-O3P-PA-O2A
2	D	601	FAD	P-O3P-PA-O1A

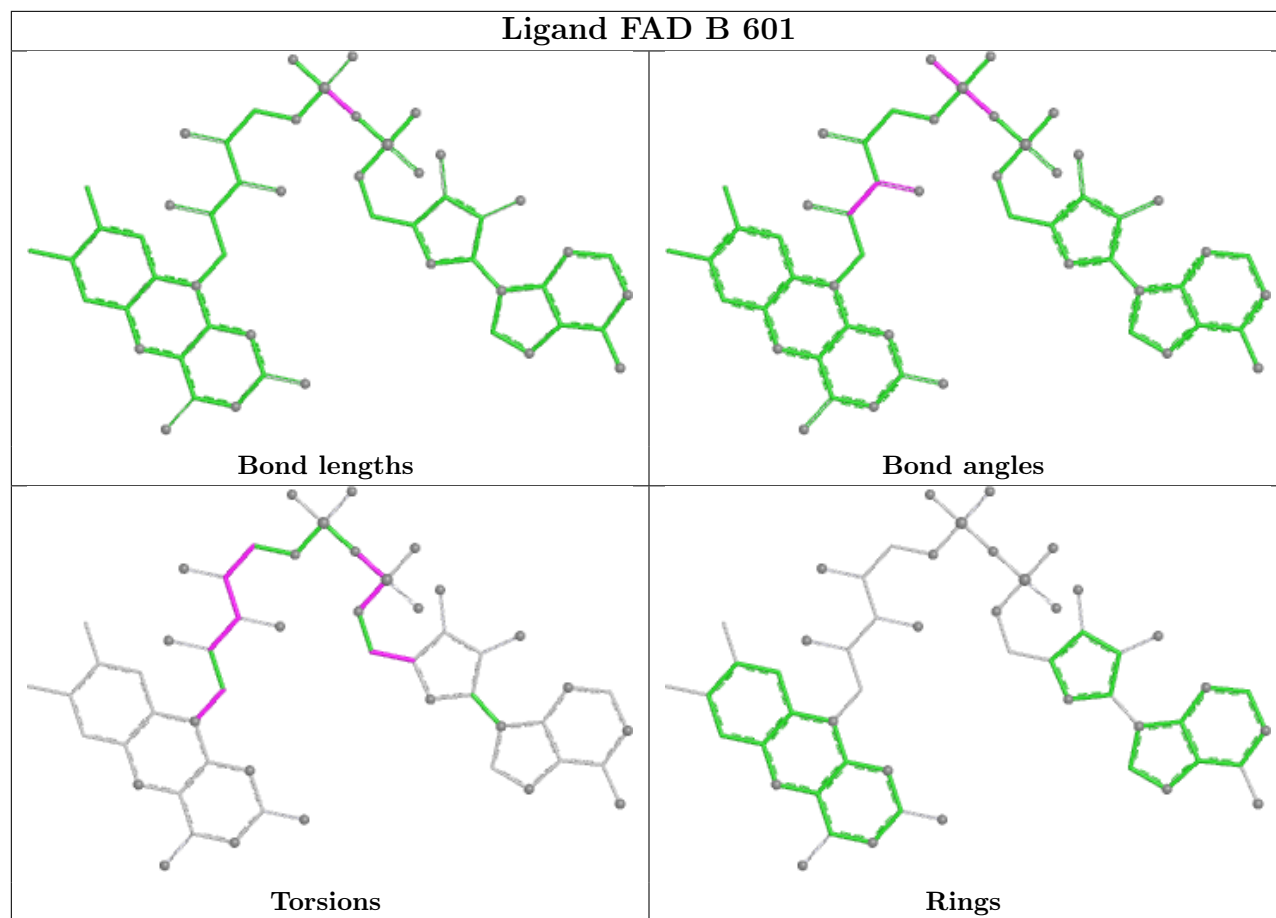
There are no ring outliers.

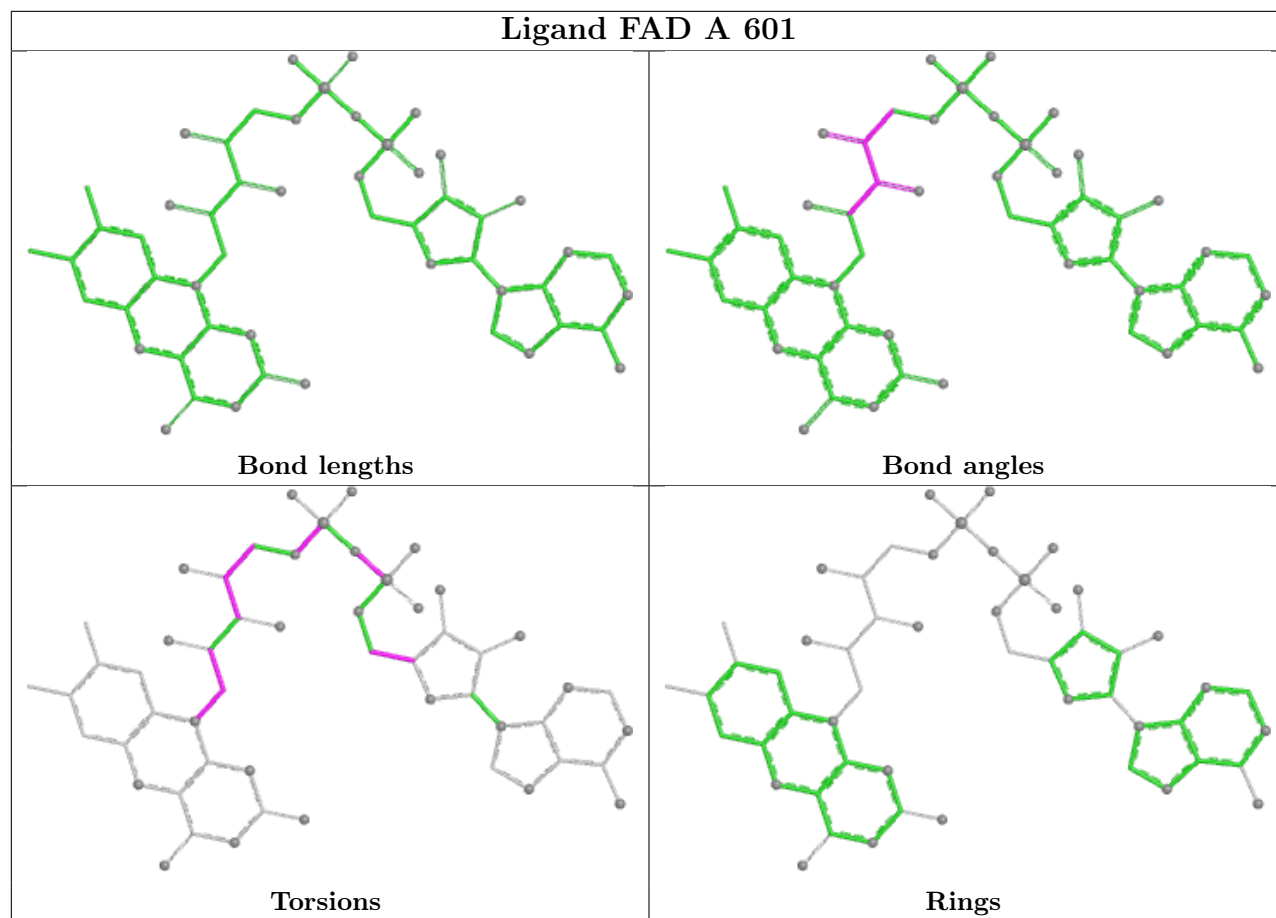
8 monomers are involved in 30 short contacts:

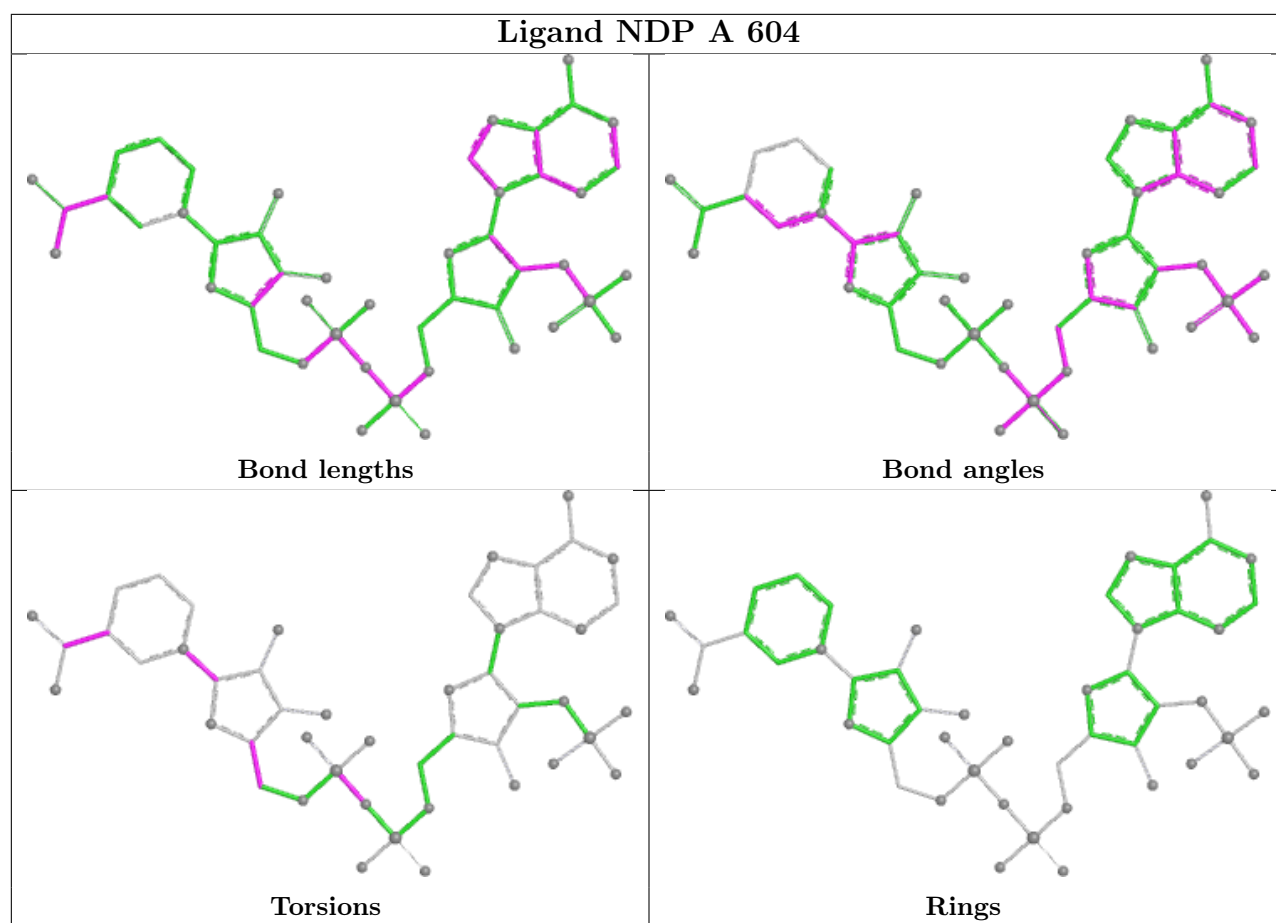
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	605	SO4	1	0
3	D	603	SO4	1	0
2	B	601	FAD	5	0
2	A	601	FAD	7	0
4	A	604	NDP	2	0
2	C	601	FAD	6	0
2	D	601	FAD	7	0
4	B	603	NDP	1	0

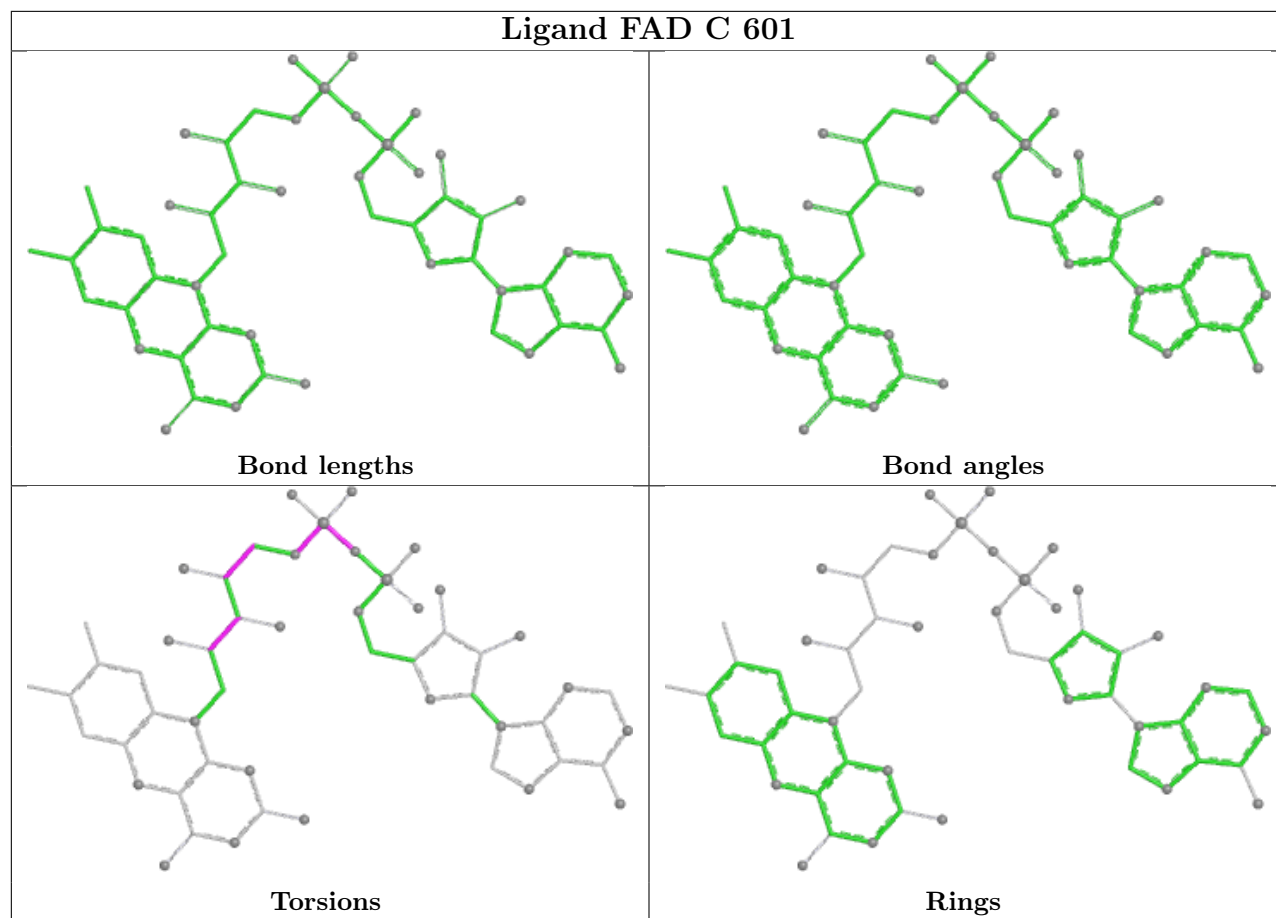
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

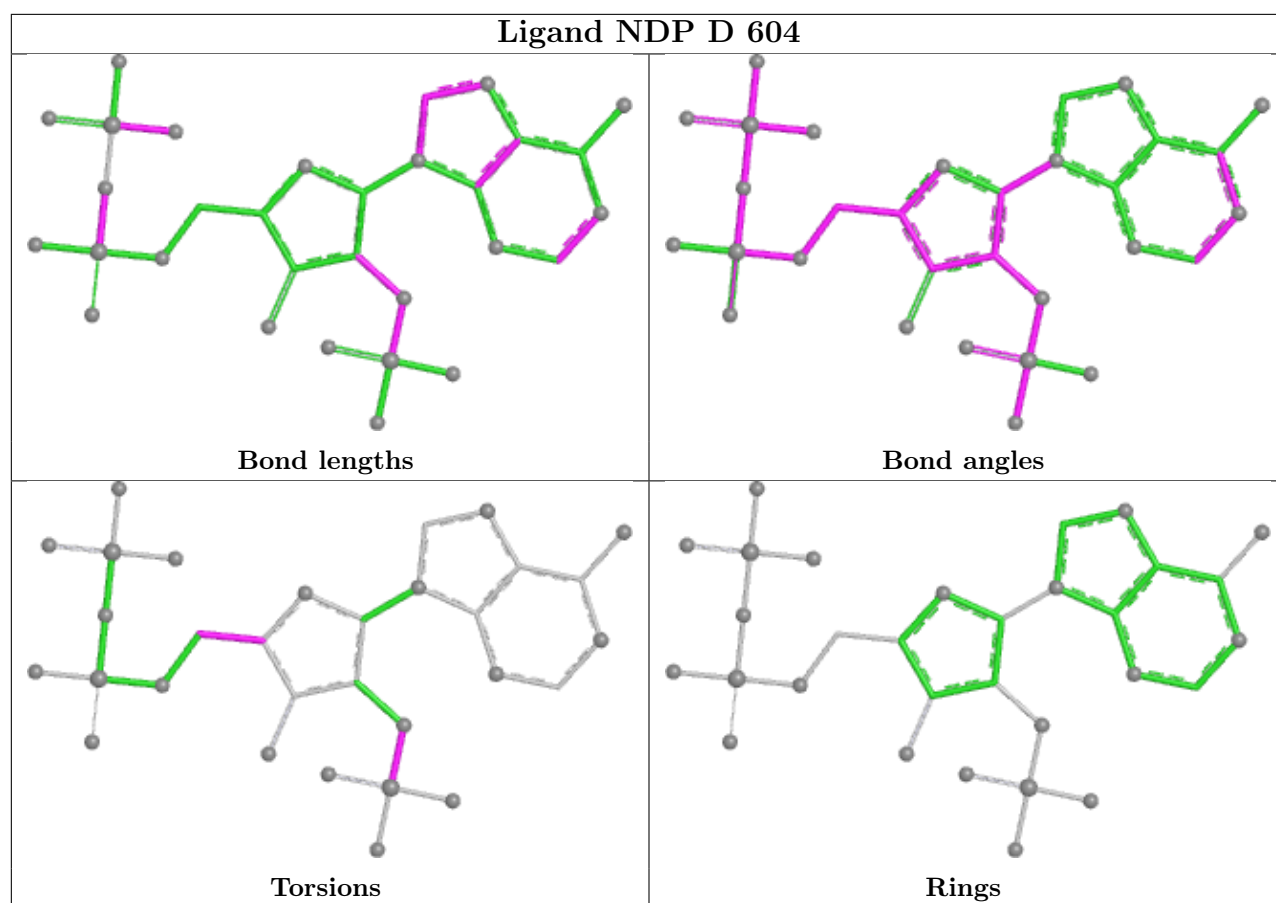


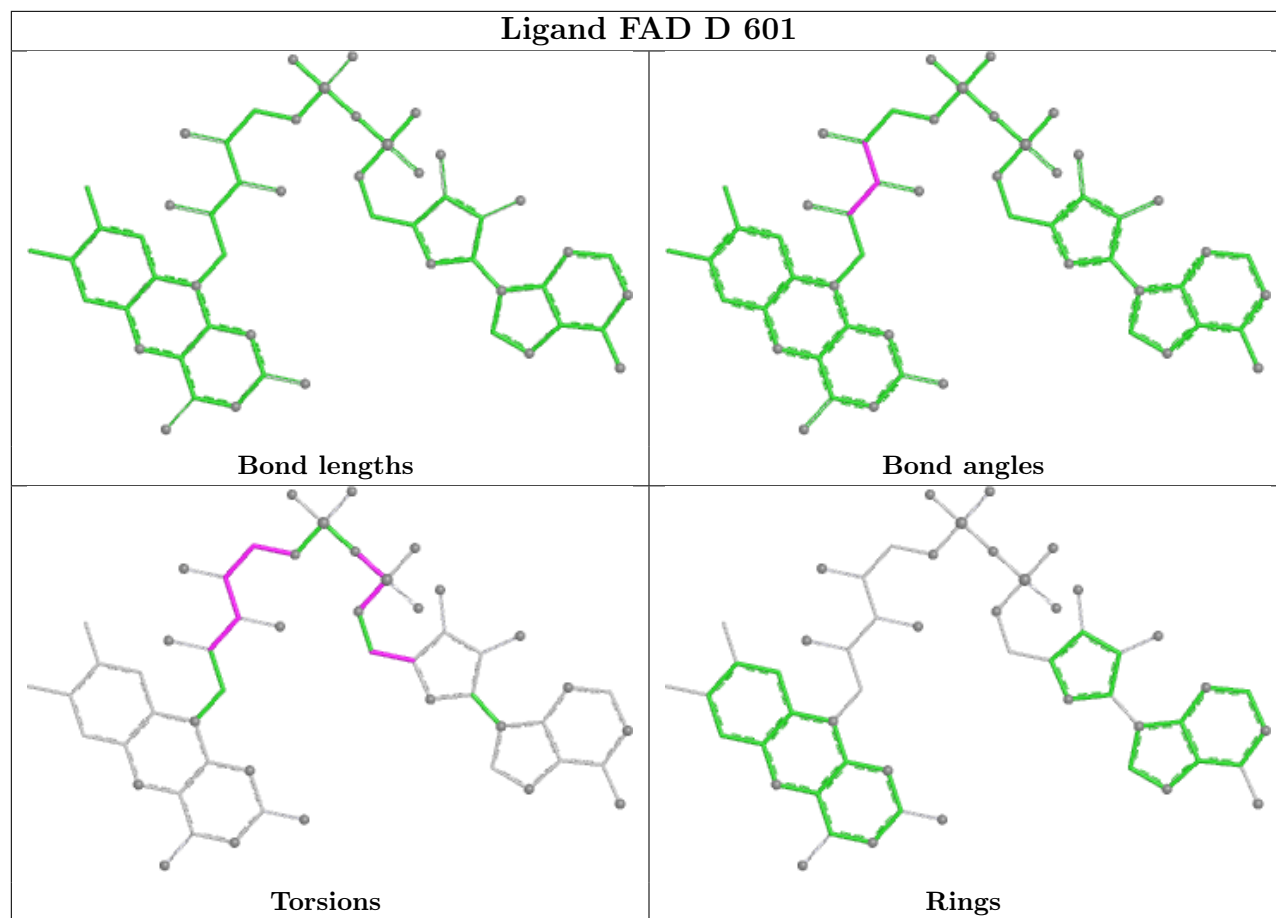


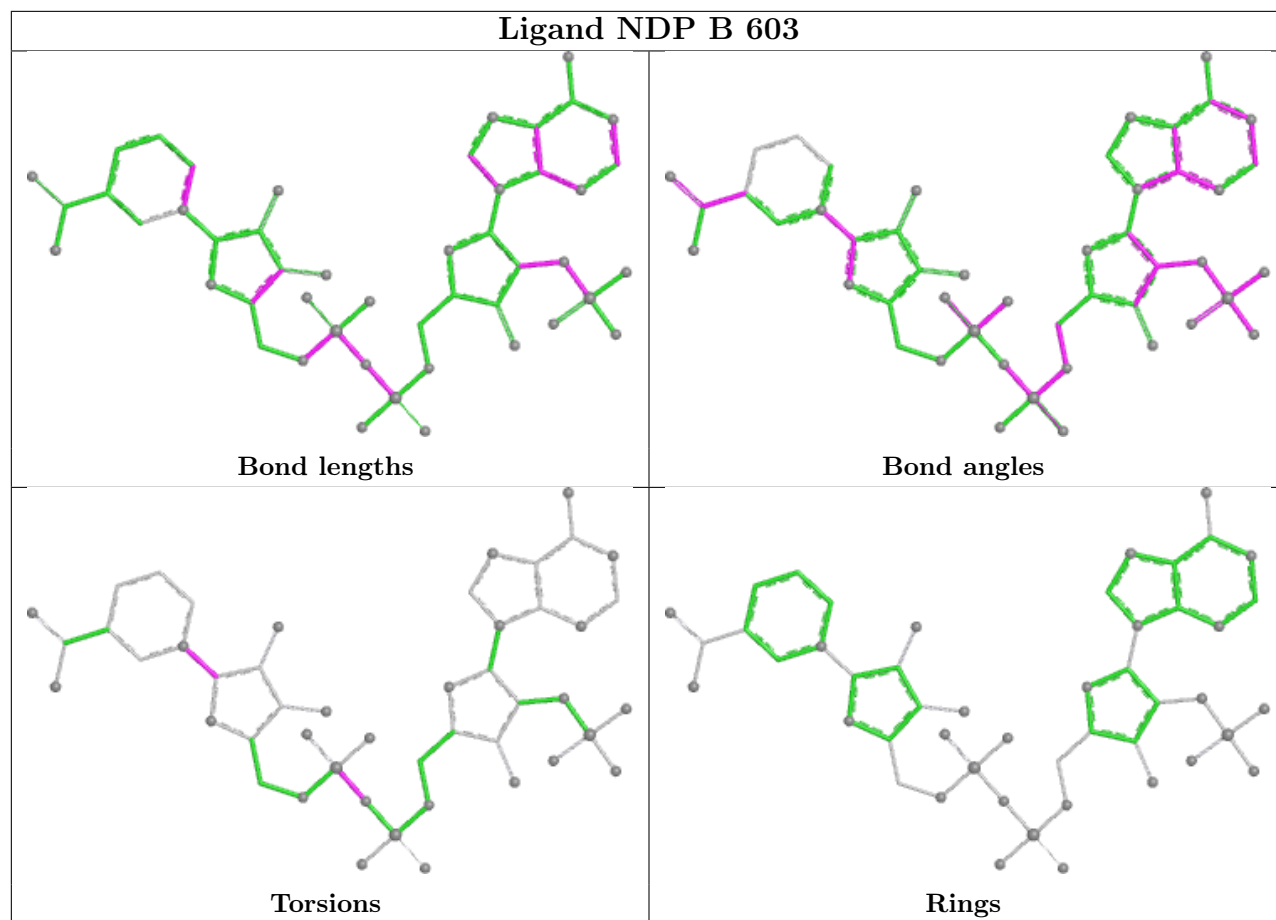












5.7 Other polymers [i](#)

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	495/513 (96%)	-0.19	11 (2%) 62 60	28, 44, 74, 101	0
1	B	492/513 (95%)	-0.18	13 (2%) 57 55	28, 47, 78, 106	0
1	C	501/513 (97%)	-0.03	24 (4%) 35 35	29, 55, 82, 102	0
1	D	499/513 (97%)	0.07	21 (4%) 40 39	27, 56, 91, 107	0
All	All	1987/2052 (96%)	-0.09	69 (3%) 47 44	27, 50, 83, 107	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	207	ASN	7.8
1	D	3	HIS	5.5
1	C	506	LYS	5.2
1	A	3	HIS	4.9
1	D	506	LYS	4.5
1	B	358	GLN	4.4
1	A	309	ARG	4.1
1	A	199	LYS	4.1
1	B	505	ALA	4.0
1	C	64	VAL	4.0
1	B	3	HIS	3.9
1	C	3	HIS	3.8
1	D	238	GLU	3.6
1	B	308	GLU	3.5
1	D	65	ILE	3.4
1	A	307	PRO	3.4
1	A	505	ALA	3.4
1	A	310	ILE	3.3
1	D	208	ALA	3.2
1	D	64	VAL	3.1
1	C	206	PRO	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	255	GLN	3.0
1	D	253	THR	3.0
1	A	6	ILE	3.0
1	C	491	GLN	3.0
1	C	248	ASN	2.9
1	D	204	TRP	2.8
1	D	202	GLY	2.8
1	C	379	MET	2.8
1	D	35	MET	2.8
1	B	260	ILE	2.7
1	C	202	GLY	2.7
1	C	249	ASN	2.7
1	A	359	SER	2.7
1	C	183	VAL	2.7
1	C	60	VAL	2.6
1	C	6	ILE	2.5
1	C	92	ILE	2.5
1	D	6	ILE	2.5
1	C	10	VAL	2.5
1	C	65	ILE	2.4
1	C	203	ASN	2.4
1	D	203	ASN	2.4
1	C	83	GLU	2.4
1	B	7	SER	2.4
1	D	442	ILE	2.4
1	C	279	ASN	2.3
1	D	4	PRO	2.3
1	C	238	GLU	2.3
1	B	208	ALA	2.2
1	A	41	GLU	2.2
1	C	505	ALA	2.2
1	C	200	THR	2.2
1	D	233	LYS	2.2
1	D	380	LYS	2.2
1	C	260	ILE	2.2
1	B	252	VAL	2.2
1	C	204	TRP	2.1
1	D	182	ARG	2.1
1	D	252	VAL	2.1
1	B	238	GLU	2.1
1	B	185	ALA	2.1
1	D	60	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	439	ASP	2.0
1	A	181	GLU	2.0
1	B	4	PRO	2.0
1	B	249	ASN	2.0
1	B	356	ARG	2.0
1	C	305	SER	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(Å ²)	Q<0.9
3	SO4	A	603	5/5	0.77	0.15	78,80,86,90	5
3	SO4	B	602	5/5	0.79	0.15	62,74,81,87	5
3	SO4	B	605	5/5	0.83	0.15	76,80,81,82	5
4	NDP	D	604	31/48	0.87	0.17	42,68,87,95	31
3	SO4	C	605	5/5	0.88	0.14	69,73,79,86	5
4	NDP	C	602	31/48	0.90	0.14	38,58,69,72	31
3	SO4	C	603	5/5	0.92	0.13	58,63,71,80	5
3	SO4	D	603	5/5	0.92	0.12	69,73,76,78	5
3	SO4	D	602	5/5	0.93	0.11	64,64,69,71	5
2	FAD	D	601	53/53	0.93	0.11	40,60,74,91	0
2	FAD	C	601	53/53	0.94	0.11	28,60,73,86	0
3	SO4	A	602	5/5	0.94	0.10	51,61,72,74	5
3	SO4	B	604	5/5	0.94	0.16	68,71,72,74	5
3	SO4	C	606	5/5	0.94	0.15	76,78,80,81	5
3	SO4	C	604	5/5	0.95	0.13	65,77,83,86	5
3	SO4	D	605	5/5	0.96	0.11	59,66,73,74	5
4	NDP	A	604	48/48	0.96	0.08	41,56,70,74	0

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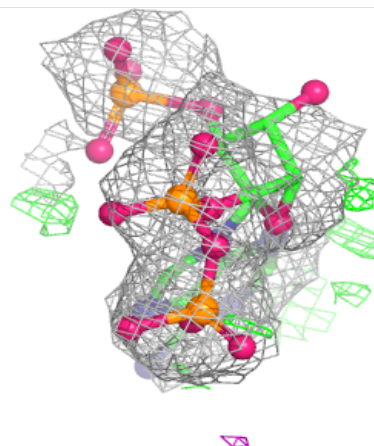
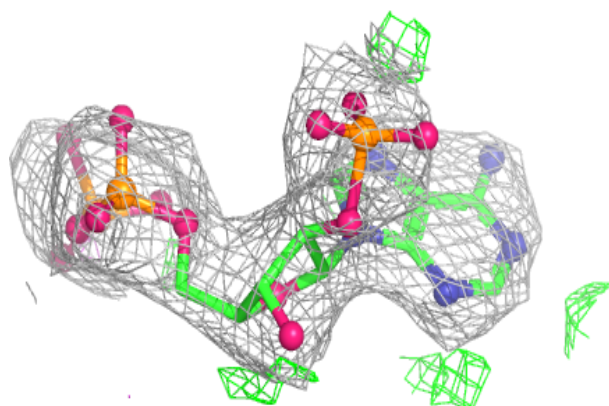
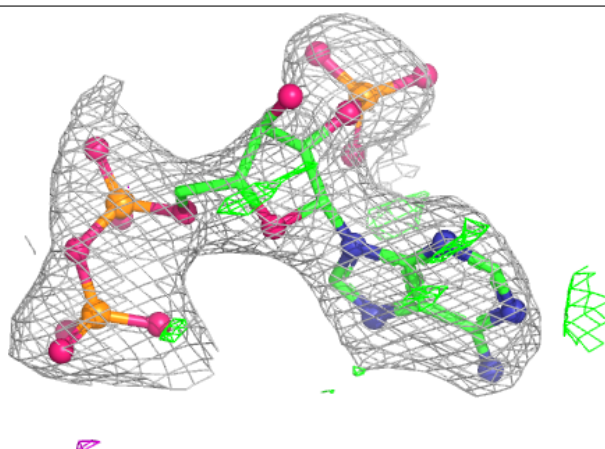
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	A	605	5/5	0.96	0.13	67,75,79,82	0
3	SO4	B	606	5/5	0.96	0.19	71,72,74,76	0
4	NDP	B	603	48/48	0.97	0.07	35,48,60,68	0
2	FAD	B	601	53/53	0.97	0.07	23,36,47,50	0
2	FAD	A	601	53/53	0.97	0.08	17,37,50,63	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.

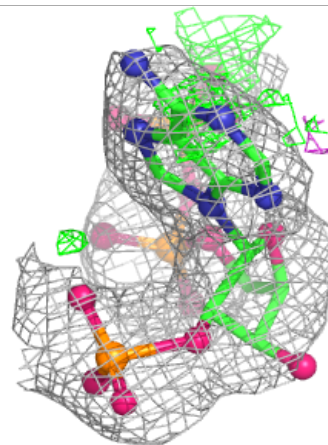
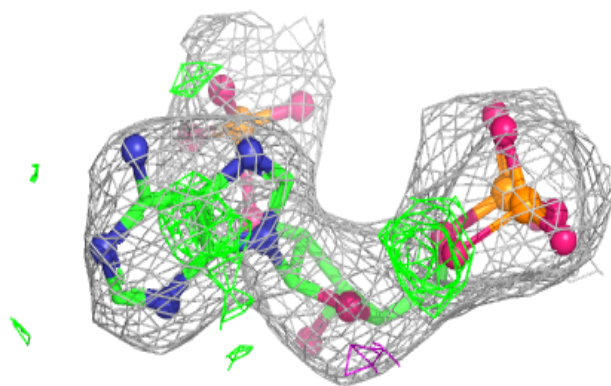
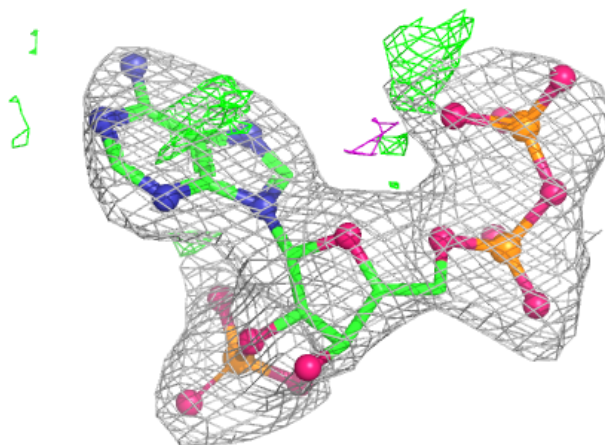
Electron density around NDP D 604:

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



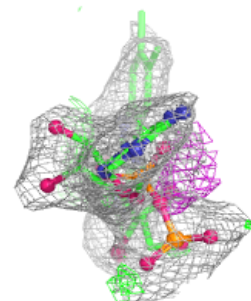
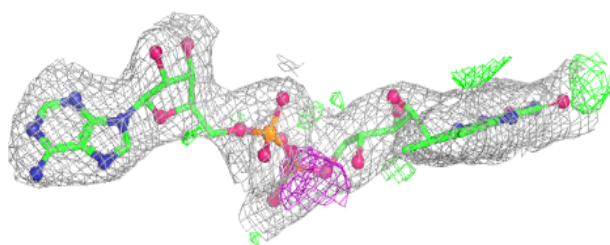
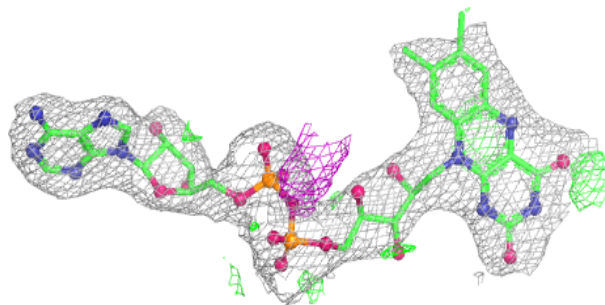
Electron density around NDP C 602:

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

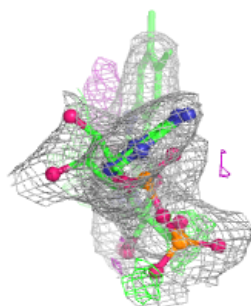
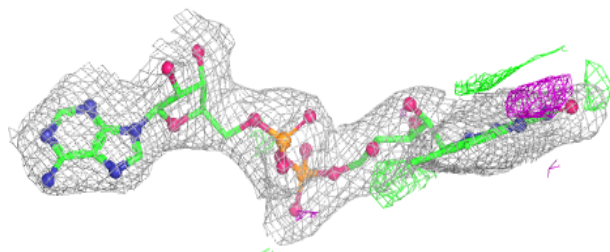
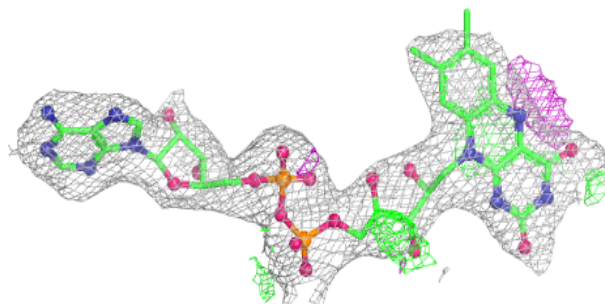


Electron density around FAD D 601:

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

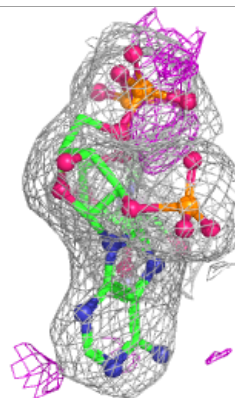
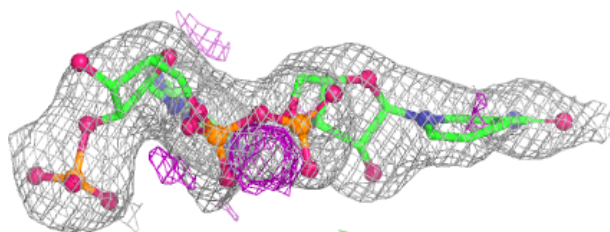
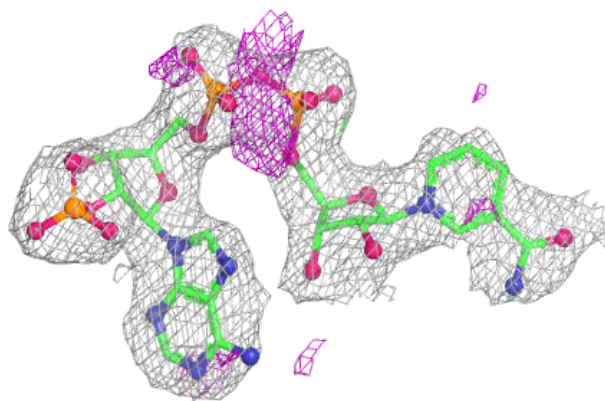
**Electron density around FAD C 601:**

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

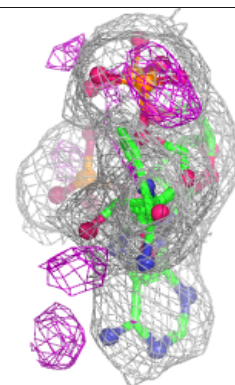
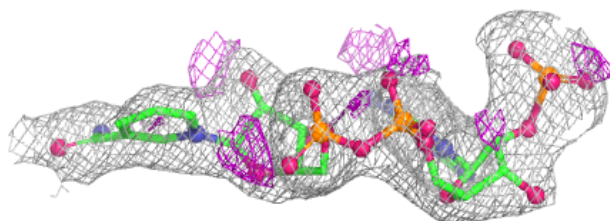
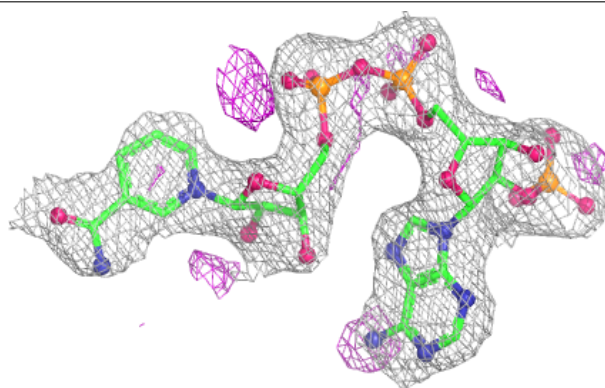


Electron density around NDP A 604:

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

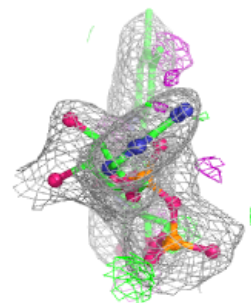
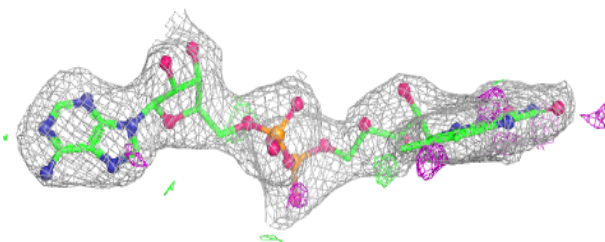
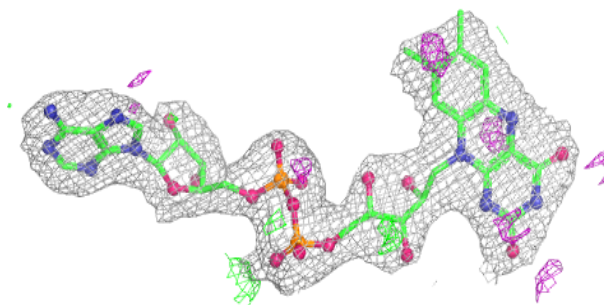
**Electron density around NDP B 603:**

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

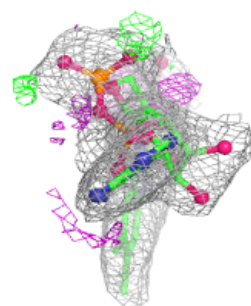
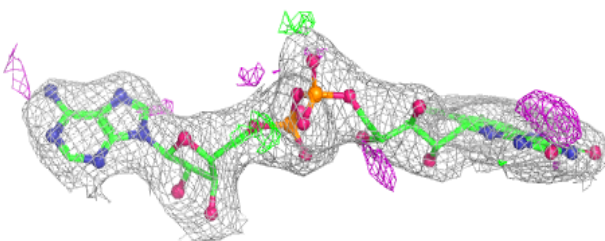
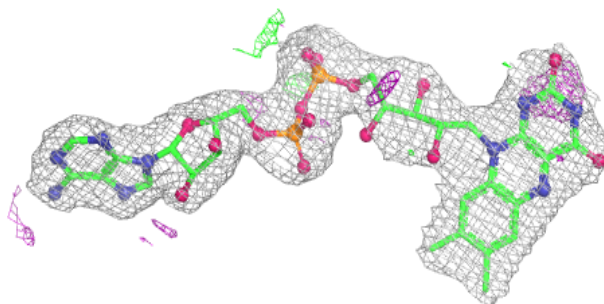


Electron density around FAD B 601:

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

**Electron density around FAD 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.