



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 11:50 AM GMT

PDB ID : 1KBQ
Title : Complex of Human NAD(P)H quinone Oxidoreductase with 5-methoxy-1,2-dimethyl-3-(4-nitrophenoxy-methyl)indole-4,7-dione(ES936)
Authors : Faig, M.; Bianchet, M.A.; Amzel, L.M.
Deposited on : 2001-11-06
Resolution : 1.80 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

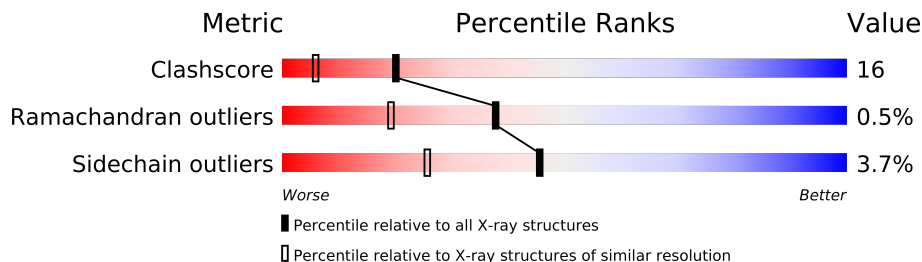
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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))
Clashscore	79885	4461 (1.80-1.80)
Ramachandran outliers	78287	4404 (1.80-1.80)
Sidechain outliers	78261	4403 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	273	
1	B	273	
1	C	273	
1	D	273	

2 Entry composition i

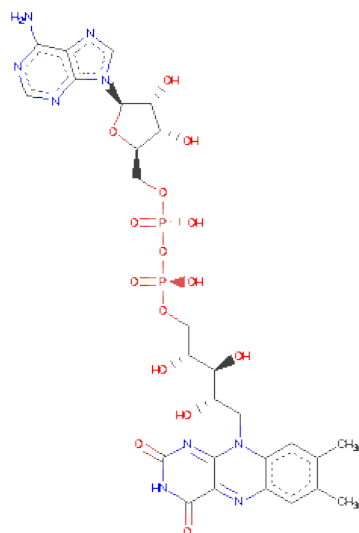
There are 4 unique types of molecules in this entry. The entry contains 9494 atoms, of which 0 are hydrogen and 0 are deuterium.

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 NAD(P)H dehydrogenase [quinone] 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	272	Total	C	N	O	S	0	0	0
			2168	1409	364	388	7			
1	B	271	Total	C	N	O	S	0	0	0
			2164	1407	363	387	7			
1	C	271	Total	C	N	O	S	0	0	0
			2164	1407	363	387	7			
1	D	272	Total	C	N	O	S	0	0	0
			2168	1409	364	388	7			

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



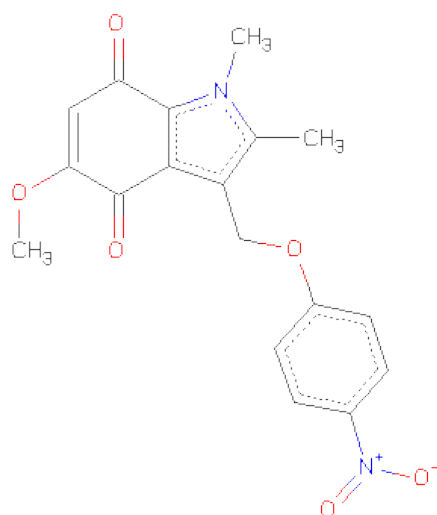
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
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 5-METHOXY-1,2-DIMETHYL-3-(4-NITROPHENOXYMETHYL)INDOLE-4,7-DIONE (three-letter code: 936) (formula: C₁₈H₁₆N₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	0
			26	18	2	6		
3	D	1	Total	C	N	O	0	0
			26	18	2	6		
3	A	1	Total	C	N	O	0	0
			26	18	2	6		
3	B	1	Total	C	N	O	0	0
			26	18	2	6		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	129	Total	O	0	0
			129	129		
4	B	144	Total	O	0	0
			144	144		
4	C	114	Total	O	0	0
			114	114		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	127	Total 127	O 127	0	0

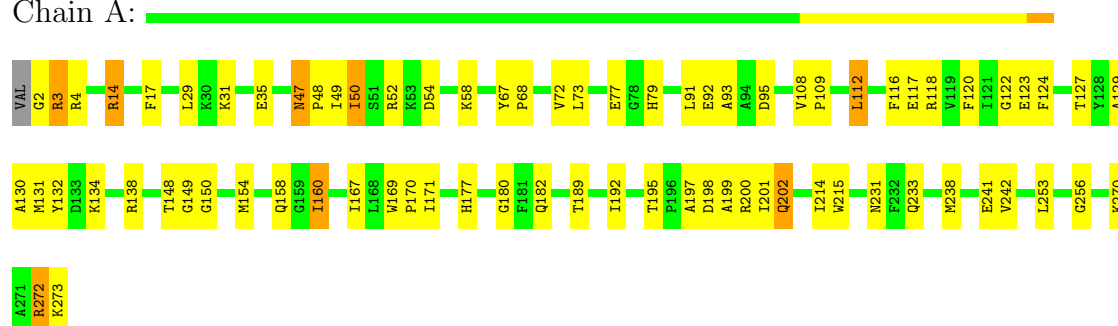
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

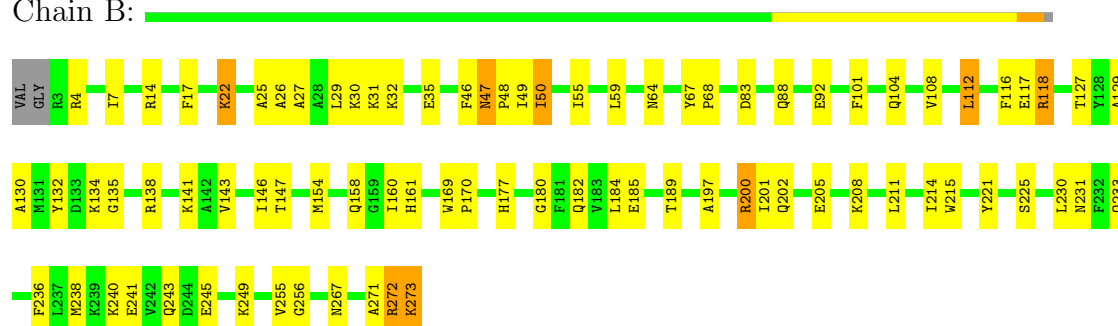
- Molecule 1: NAD(P)H dehydrogenase [quinone] 1

Chain A:



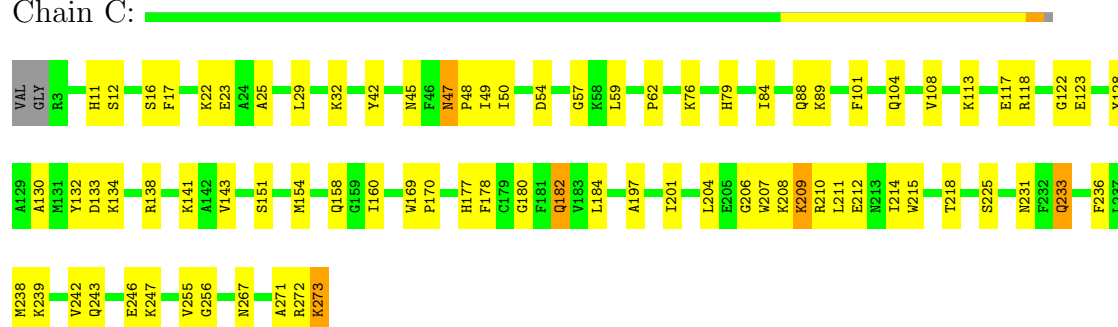
- Molecule 1: NAD(P)H dehydrogenase [quinone] 1

Chain B:



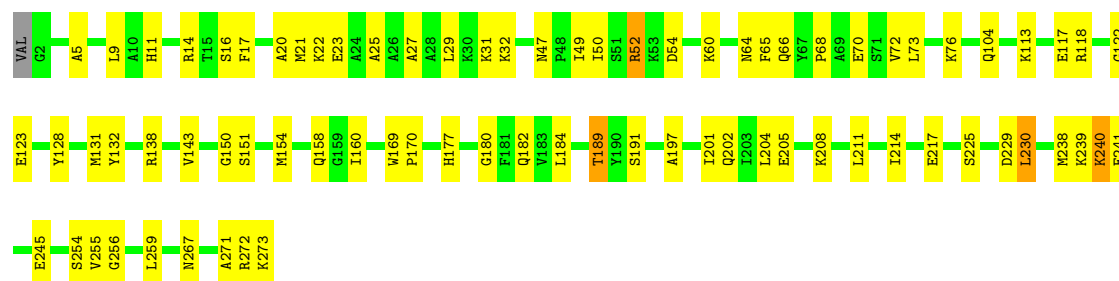
- Molecule 1: NAD(P)H dehydrogenase [quinone] 1

Chain C:



- Molecule 1: NAD(P)H dehydrogenase [quinone] 1

Chain D: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.48Å 56.77Å 96.91Å 77.25° 76.83° 86.90°	Depositor
Resolution (Å)	32.19 – 1.80	Depositor
% Data completeness (in resolution range)	94.6 (32.19-1.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.215 , 0.252	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9494	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 936, FAD

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	0/2226	0.64	1/3005 (0.0%)
1	B	0.41	0/2222	0.66	3/3000 (0.1%)
1	C	0.39	0/2222	0.60	1/3000 (0.0%)
1	D	0.41	0/2226	0.61	1/3005 (0.0%)
All	All	0.40	0/8896	0.63	6/12010 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	118	ARG	NE-CZ-NH1	6.05	123.33	120.30
1	D	182	GLN	N-CA-C	-6.05	94.66	111.00
1	A	182	GLN	N-CA-C	-5.83	95.27	111.00
1	C	182	GLN	N-CA-C	-5.75	95.49	111.00
1	B	118	ARG	NE-CZ-NH2	-5.71	117.44	120.30
1	B	182	GLN	N-CA-C	-5.13	97.15	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2168	0	2167	71	0
1	B	2164	0	2164	77	0
1	C	2164	0	2164	65	0
1	D	2168	0	2167	82	0
2	A	53	0	31	6	0
2	B	53	0	31	3	0
2	C	53	0	31	3	0
2	D	53	0	31	3	0
3	A	26	0	16	1	0
3	B	26	0	16	2	0
3	C	26	0	16	2	0
3	D	26	0	16	3	0
4	A	129	0	0	5	0
4	B	144	0	0	6	0
4	C	114	0	0	3	0
4	D	127	0	0	5	0
All	All	9494	0	8850	284	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 16.

All (284) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:108:VAL:CG2	1:B:112:LEU:HD13	1.94	0.97
1:B:135:GLY:O	1:B:138:ARG:HG3	1.68	0.93
1:A:14:ARG:H	1:A:14:ARG:NE	1.68	0.92
1:D:17:PHE:HB2	2:D:604:FAD:H52A	1.51	0.92
1:A:50:ILE:HG13	1:A:118:ARG:HG2	1.53	0.90
1:A:14:ARG:HE	1:A:14:ARG:H	0.90	0.90
1:D:64:ASN:HD21	1:D:66:GLN:HE21	1.21	0.86
1:D:143:VAL:HG12	1:D:184:LEU:HB2	1.55	0.86
1:A:50:ILE:HD12	1:A:117:GLU:HB3	1.57	0.86
1:A:17:PHE:HB2	2:A:601:FAD:H52A	1.60	0.83
1:D:143:VAL:CG1	1:D:184:LEU:HB2	2.08	0.83
1:B:108:VAL:HG21	1:B:112:LEU:HD13	1.61	0.82
1:A:108:VAL:CG1	1:A:112:LEU:HB3	2.11	0.81
1:B:48:PRO:HG3	1:D:49:ILE:HD11	1.64	0.79
1:C:206:GLY:O	1:C:209:LYS:HD3	1.82	0.79
1:A:14:ARG:HE	1:A:14:ARG:N	1.75	0.78
1:B:17:PHE:HB2	2:B:602:FAD:H52A	1.65	0.77
1:D:64:ASN:HD21	1:D:66:GLN:NE2	1.83	0.77
1:C:50:ILE:HG22	1:C:118:ARG:HG2	1.66	0.77
1:D:60:LYS:HG3	1:D:70:GLU:OE1	1.84	0.76

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:143:VAL:HG22	1:B:184:LEU:HB2	1.68	0.74
1:C:17:PHE:HB2	2:C:603:FAD:H52A	1.69	0.73
1:A:127:THR:HG22	1:A:129:ALA:H	1.54	0.73
1:A:68:PRO:O	1:A:72:VAL:HG23	1.88	0.73
1:A:58:LYS:N	1:A:58:LYS:HD2	2.02	0.73
1:D:64:ASN:ND2	1:D:66:GLN:HE21	1.88	0.72
1:C:151:SER:H	1:C:154:MET:HE3	1.54	0.72
1:D:60:LYS:HG2	1:D:73:LEU:CD2	2.18	0.72
1:B:231:ASN:ND2	1:B:233:GLN:HB3	2.05	0.72
1:C:273:LYS:HD2	1:C:273:LYS:H	1.54	0.72
1:A:273:LYS:HG2	1:A:273:LYS:OXT	1.91	0.71
1:B:236:PHE:HB3	1:D:160:ILE:HG13	1.74	0.69
1:A:50:ILE:CD1	1:A:117:GLU:HB3	2.22	0.69
1:B:160:ILE:HD12	1:D:131:MET:HE2	1.74	0.69
1:A:189:THR:HG22	4:A:777:HOH:O	1.92	0.69
1:D:151:SER:H	1:D:154:MET:HE3	1.55	0.69
1:A:148:THR:HG23	2:A:601:FAD:O2	1.93	0.69
1:D:21:MET:CE	1:D:204:LEU:HD23	2.23	0.69
1:B:231:ASN:HD21	1:B:233:GLN:HB3	1.58	0.69
1:D:189:THR:HG23	4:D:719:HOH:O	1.94	0.68
1:D:52:ARG:HG2	4:D:752:HOH:O	1.94	0.67
1:A:160:ILE:HG13	1:C:236:PHE:HB3	1.76	0.67
1:C:89:LYS:HG3	4:C:748:HOH:O	1.95	0.66
1:D:255:VAL:HG23	1:D:267:ASN:HD22	1.60	0.66
1:D:169:TRP:CZ2	1:D:256:GLY:HA3	2.31	0.66
1:A:160:ILE:CG1	1:C:236:PHE:HB3	2.27	0.65
1:C:57:GLY:H	1:C:79:HIS:HB3	1.60	0.65
1:B:160:ILE:HD12	1:D:131:MET:CE	2.26	0.65
1:D:25:ALA:O	1:D:29:LEU:HD13	1.96	0.65
1:C:238:MET:HE3	1:C:243:GLN:HG2	1.79	0.65
1:D:273:LYS:HB3	4:D:750:HOH:O	1.96	0.64
1:C:204:LEU:O	1:C:208:LYS:HG3	1.97	0.64
1:A:189:THR:HG23	4:A:773:HOH:O	1.98	0.64
1:B:238:MET:HE3	1:B:243:GLN:HG2	1.80	0.64
1:C:59:LEU:HB2	1:C:62:PRO:HG3	1.80	0.63
1:C:243:GLN:O	1:C:247:LYS:HG3	1.98	0.63
1:A:200:ARG:NH1	2:A:601:FAD:H1B	2.13	0.63
1:D:189:THR:HG22	4:D:773:HOH:O	1.98	0.63
1:B:88:GLN:O	1:B:92:GLU:HG3	1.99	0.63
1:D:50:ILE:HG22	1:D:118:ARG:HG2	1.80	0.63
1:B:50:ILE:HG12	1:B:67:TYR:CZ	2.34	0.63
1:D:60:LYS:HG2	1:D:73:LEU:HD22	1.80	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:50:ILE:HD12	1:B:117:GLU:HB3	1.80	0.62
1:A:108:VAL:HG11	1:A:112:LEU:HB3	1.81	0.62
1:B:25:ALA:O	1:B:29:LEU:HD13	2.00	0.61
1:D:21:MET:HE2	1:D:204:LEU:HD23	1.83	0.61
1:B:154:MET:HE3	1:B:160:ILE:HD11	1.83	0.61
1:D:31:LYS:HB3	1:D:31:LYS:NZ	2.15	0.61
1:D:60:LYS:CD	1:D:73:LEU:HD22	2.31	0.60
1:B:112:LEU:HD22	1:B:116:PHE:CE2	2.37	0.60
1:B:134:LYS:HA	1:B:138:ARG:HD3	1.82	0.60
1:A:108:VAL:CG1	1:A:112:LEU:HD13	2.31	0.60
1:A:198:ASP:O	1:A:202:GLN:HG2	2.02	0.60
1:D:76:LYS:HE2	1:D:123:GLU:OE2	2.02	0.60
1:B:27:ALA:O	1:B:31:LYS:HG3	2.01	0.60
1:D:201:ILE:O	1:D:205:GLU:HG2	2.02	0.60
1:A:3:ARG:HB3	1:A:95:ASP:OD2	2.02	0.59
1:A:2:GLY:O	1:A:3:ARG:HB2	2.03	0.59
1:A:108:VAL:HG11	1:A:112:LEU:HD13	1.83	0.59
1:C:169:TRP:CZ2	1:C:256:GLY:HA3	2.38	0.59
1:A:270:LYS:HB2	1:A:272:ARG:NH1	2.18	0.59
1:A:50:ILE:HD11	1:A:117:GLU:O	2.03	0.58
1:B:134:LYS:HE3	1:B:225:SER:OG	2.03	0.58
1:D:138:ARG:HA	1:D:180:GLY:O	2.03	0.58
1:A:4:ARG:HD3	1:A:35:GLU:OE1	2.04	0.57
1:C:169:TRP:HB3	1:C:170:PRO:HD3	1.85	0.57
1:C:130:ALA:HB1	1:C:134:LYS:O	2.04	0.57
1:A:197:ALA:O	1:A:201:ILE:HG12	2.04	0.57
1:C:122:GLY:O	1:C:123:GLU:HB2	2.05	0.57
1:A:58:LYS:H	1:A:58:LYS:HD2	1.68	0.57
1:B:108:VAL:HG22	1:B:112:LEU:HB3	1.87	0.56
1:D:27:ALA:O	1:D:31:LYS:HG3	2.06	0.56
1:A:132:TYR:O	1:A:180:GLY:HA2	2.05	0.56
1:C:246:GLU:HG3	4:C:733:HOH:O	2.06	0.56
1:B:271:ALA:O	1:B:272:ARG:HB2	2.06	0.56
1:A:50:ILE:HG12	1:A:67:TYR:CZ	2.41	0.56
1:D:20:ALA:HA	1:D:23:GLU:OE2	2.05	0.56
1:B:4:ARG:HB3	4:B:796:HOH:O	2.06	0.56
1:D:169:TRP:HB3	1:D:170:PRO:HD3	1.88	0.55
1:B:50:ILE:HG13	1:B:118:ARG:HG2	1.87	0.55
1:D:122:GLY:O	1:D:123:GLU:HB2	2.05	0.55
1:D:197:ALA:O	1:D:201:ILE:HG13	2.06	0.55
1:B:169:TRP:HB3	1:B:170:PRO:HD3	1.86	0.55
1:B:236:PHE:HD2	1:D:160:ILE:HD12	1.72	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:230:LEU:HD22	1:D:160:ILE:HD13	1.87	0.55
1:D:21:MET:HE3	1:D:204:LEU:HD23	1.88	0.55
1:B:141:LYS:HD3	1:B:184:LEU:HD21	1.88	0.55
1:D:204:LEU:O	1:D:208:LYS:HG3	2.06	0.55
1:B:202:GLN:NE2	1:B:202:GLN:HA	2.21	0.54
1:A:231:ASN:ND2	1:A:233:GLN:HE21	2.05	0.54
1:B:29:LEU:HD11	1:B:211:LEU:HB3	1.90	0.54
1:C:25:ALA:O	1:C:29:LEU:HD13	2.07	0.54
1:C:76:LYS:HE3	1:C:123:GLU:HG3	1.89	0.54
1:A:138:ARG:HA	1:A:180:GLY:O	2.08	0.54
1:C:233:GLN:HA	1:C:233:GLN:HE21	1.72	0.54
1:A:2:GLY:O	1:A:3:ARG:CB	2.56	0.53
1:A:4:ARG:HD2	1:A:93:ALA:O	2.08	0.53
1:B:169:TRP:CZ2	1:B:256:GLY:HA3	2.44	0.53
1:B:127:THR:HG22	1:B:129:ALA:H	1.74	0.53
1:C:143:VAL:HG22	1:C:184:LEU:HB2	1.90	0.53
1:C:47:ASN:ND2	1:C:49:ILE:H	2.07	0.53
4:B:837:HOH:O	1:D:52:ARG:HD3	2.08	0.53
1:B:238:MET:CE	1:B:243:GLN:HG2	2.37	0.53
1:B:143:VAL:CG2	1:B:184:LEU:HD12	2.39	0.52
1:B:154:MET:HE1	1:B:161:HIS:NE2	2.24	0.52
1:B:132:TYR:O	1:B:180:GLY:HA2	2.09	0.52
1:D:17:PHE:HB2	2:D:604:FAD:C5B	2.33	0.52
1:B:4:ARG:HG2	1:B:35:GLU:OE1	2.09	0.52
1:D:23:GLU:HB3	4:D:784:HOH:O	2.10	0.52
1:A:200:ARG:HH11	2:A:601:FAD:H1B	1.74	0.52
1:B:134:LYS:HB2	1:B:134:LYS:NZ	2.25	0.52
1:A:91:LEU:HD11	1:A:120:PHE:HE1	1.75	0.51
1:D:60:LYS:CG	1:D:73:LEU:HD22	2.40	0.51
1:D:76:LYS:CE	1:D:123:GLU:HG3	2.41	0.51
1:A:131:MET:HE3	1:C:160:ILE:HD13	1.91	0.51
1:B:47:ASN:ND2	1:B:49:ILE:H	2.09	0.50
1:C:84:ILE:O	1:C:88:GLN:HG3	2.11	0.50
1:B:26:ALA:O	1:B:30:LYS:HG3	2.12	0.50
1:B:255:VAL:HG23	1:B:267:ASN:HD22	1.76	0.50
1:D:104:GLN:HA	2:D:604:FAD:C5X	2.42	0.50
1:C:143:VAL:CG2	1:C:184:LEU:HD12	2.41	0.50
1:B:46:PHE:O	1:B:48:PRO:HD3	2.12	0.50
1:A:108:VAL:HG12	1:A:112:LEU:HB3	1.92	0.49
1:D:131:MET:SD	3:D:702:936:H281	2.52	0.49
1:A:270:LYS:HB3	1:A:272:ARG:HD3	1.94	0.49
1:B:147:THR:HG22	1:B:189:THR:OG1	2.13	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:132:TYR:O	1:C:180:GLY:HA2	2.12	0.49
1:C:57:GLY:HA3	1:C:79:HIS:CG	2.47	0.49
1:A:169:TRP:CZ2	1:A:256:GLY:HA3	2.47	0.49
1:C:197:ALA:O	1:C:201:ILE:HG13	2.11	0.49
1:B:132:TYR:HA	1:B:177:HIS:O	2.13	0.49
1:B:214:ILE:HG23	1:B:215:TRP:N	2.29	0.48
1:B:64:ASN:ND2	4:B:807:HOH:O	2.46	0.48
1:A:148:THR:HG22	1:A:150:GLY:N	2.29	0.48
1:D:214:ILE:HD12	1:D:217:GLU:OE2	2.14	0.48
1:D:20:ALA:HA	1:D:23:GLU:HG2	1.95	0.48
1:C:242:VAL:O	1:C:246:GLU:HB2	2.14	0.48
1:B:185:GLU:HG3	4:B:714:HOH:O	2.13	0.48
1:C:113:LYS:O	1:C:117:GLU:HG3	2.13	0.48
1:B:230:LEU:HD22	1:D:160:ILE:CD1	2.43	0.48
3:C:701:936:H291	3:C:701:936:H192	1.68	0.48
1:C:151:SER:H	1:C:154:MET:CE	2.24	0.48
1:B:104:GLN:HA	2:B:602:FAD:C5X	2.44	0.47
1:B:189:THR:HG23	4:B:712:HOH:O	2.13	0.47
1:B:55:ILE:HG21	1:B:59:LEU:HD23	1.95	0.47
1:C:76:LYS:CE	1:C:123:GLU:HG3	2.45	0.47
1:C:132:TYR:CD1	1:C:178:PHE:HA	2.49	0.47
1:A:132:TYR:HA	1:A:177:HIS:O	2.14	0.47
1:C:141:LYS:HE2	1:C:182:GLN:OE1	2.14	0.47
1:A:47:ASN:ND2	1:A:49:ILE:H	2.11	0.47
1:A:192:ILE:CG2	2:A:601:FAD:H5'1	2.45	0.47
1:C:132:TYR:HA	1:C:177:HIS:O	2.14	0.47
1:D:9:LEU:HD22	1:D:22:LYS:HD3	1.96	0.47
1:A:52:ARG:HG2	4:A:719:HOH:O	2.14	0.47
1:B:236:PHE:HD2	1:D:160:ILE:CD1	2.27	0.46
1:D:241:GLU:O	1:D:245:GLU:HG3	2.15	0.46
1:D:128:TYR:HB3	3:D:702:936:O35	2.14	0.46
1:A:231:ASN:HD21	1:A:233:GLN:HE21	1.63	0.46
1:A:73:LEU:HD23	4:A:780:HOH:O	2.16	0.46
1:D:255:VAL:H	1:D:267:ASN:ND2	2.14	0.46
1:D:254:SER:HB2	1:D:267:ASN:HD21	1.80	0.46
1:B:67:TYR:HB3	1:B:68:PRO:HD3	1.98	0.46
1:C:207:TRP:HZ3	1:C:210:ARG:HD3	1.80	0.46
1:A:253:LEU:HD22	4:A:793:HOH:O	2.15	0.46
1:C:104:GLN:HA	2:C:603:FAD:C5X	2.46	0.46
1:C:12:SER:HB3	1:C:42:TYR:CE1	2.51	0.46
3:D:702:936:H192	3:D:702:936:H291	1.56	0.45
1:A:231:ASN:HD21	1:A:233:GLN:NE2	2.14	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:29:LEU:CD1	1:C:211:LEU:HD13	2.46	0.45
1:B:31:LYS:NZ	1:B:208:LYS:NZ	2.64	0.45
1:A:195:THR:HG22	1:A:199:ALA:HB3	1.98	0.45
1:C:218:THR:HG23	1:C:271:ALA:HB3	1.97	0.45
1:A:130:ALA:HB1	1:A:134:LYS:O	2.16	0.45
1:B:108:VAL:CG2	1:B:112:LEU:CD1	2.81	0.45
1:B:29:LEU:CD1	1:B:211:LEU:HB3	2.47	0.45
1:C:255:VAL:HG23	1:C:267:ASN:HD22	1.82	0.45
1:B:197:ALA:HA	1:B:200:ARG:NH1	2.32	0.45
1:B:32:LYS:HD2	4:B:785:HOH:O	2.17	0.45
1:B:83:ASP:OD2	1:B:118:ARG:NH2	2.33	0.45
1:C:23:GLU:OE1	1:C:23:GLU:HA	2.16	0.45
1:C:45:ASN:HB3	4:C:772:HOH:O	2.16	0.45
1:D:238:MET:HE2	1:D:259:LEU:HD21	1.99	0.45
1:D:132:TYR:O	1:D:180:GLY:HA2	2.16	0.45
1:A:108:VAL:O	1:C:113:LYS:NZ	2.50	0.45
1:D:60:LYS:HG2	1:D:73:LEU:HD23	1.93	0.45
1:B:135:GLY:O	1:B:138:ARG:NH1	2.49	0.44
1:A:67:TYR:HB3	1:A:68:PRO:HD3	1.99	0.44
1:C:50:ILE:CG2	1:C:118:ARG:HG2	2.44	0.44
1:B:160:ILE:CD1	1:D:131:MET:HE2	2.43	0.44
1:D:21:MET:HA	1:D:21:MET:HE2	1.99	0.44
1:B:108:VAL:HG23	1:B:112:LEU:HD13	1.94	0.44
1:A:122:GLY:O	1:A:123:GLU:HB3	2.17	0.44
1:C:32:LYS:HD2	1:C:212:GLU:HG2	1.98	0.44
3:B:704:936:H291	3:B:704:936:H192	1.80	0.44
1:D:132:TYR:HA	1:D:177:HIS:O	2.18	0.44
1:C:239:LYS:HB2	1:C:242:VAL:HG23	1.99	0.44
1:D:52:ARG:HG2	1:D:52:ARG:H	1.53	0.44
1:A:4:ARG:HD3	1:A:35:GLU:CD	2.37	0.44
1:C:273:LYS:H	1:C:273:LYS:CD	2.27	0.44
1:C:231:ASN:OD1	1:C:233:GLN:HB3	2.17	0.44
1:D:60:LYS:HD3	1:D:73:LEU:HD22	1.99	0.43
1:D:211:LEU:HA	1:D:214:ILE:HB	2.01	0.43
1:A:108:VAL:HG23	1:A:171:ILE:HD13	2.00	0.43
1:B:201:ILE:O	1:B:205:GLU:HG2	2.18	0.43
1:D:31:LYS:HB3	1:D:31:LYS:HZ3	1.83	0.43
1:D:76:LYS:HE2	1:D:123:GLU:HG3	1.99	0.43
1:A:77:GLU:HB2	1:A:79:HIS:CE1	2.54	0.43
1:D:113:LYS:O	1:D:117:GLU:HG3	2.19	0.43
1:D:47:ASN:ND2	1:D:49:ILE:H	2.16	0.43
1:D:11:HIS:CE1	1:D:16:SER:HB3	2.54	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:50:ILE:HD11	1:A:117:GLU:C	2.39	0.43
1:A:73:LEU:O	1:A:77:GLU:HG3	2.19	0.43
3:B:704:936:H251	1:D:150:GLY:HA2	2.01	0.43
1:D:271:ALA:O	1:D:272:ARG:HB2	2.19	0.43
1:C:233:GLN:CA	1:C:233:GLN:HE21	2.31	0.43
1:D:65:PHE:CE2	1:D:70:GLU:HG3	2.53	0.42
1:C:22:LYS:HE3	1:C:23:GLU:OE1	2.19	0.42
1:B:127:THR:HB	1:B:130:ALA:HB3	2.01	0.42
1:D:32:LYS:NZ	1:D:32:LYS:HB3	2.34	0.42
1:C:17:PHE:HB2	2:C:603:FAD:C5B	2.46	0.42
1:D:5:ALA:CB	1:D:29:LEU:HD23	2.50	0.42
1:D:29:LEU:CD1	1:D:211:LEU:HD13	2.49	0.42
1:B:230:LEU:CD2	1:D:160:ILE:HD13	2.50	0.42
1:A:31:LYS:HE3	1:A:31:LYS:HB2	1.79	0.42
1:A:238:MET:CE	1:A:242:VAL:HG12	2.49	0.42
1:A:148:THR:CG2	1:A:149:GLY:N	2.82	0.42
1:C:206:GLY:HA2	1:C:209:LYS:HD2	2.02	0.42
1:D:238:MET:HE3	1:D:238:MET:HB2	1.76	0.42
1:A:17:PHE:HB2	2:A:601:FAD:C5B	2.42	0.41
1:A:154:MET:HE2	1:A:160:ILE:CD1	2.50	0.41
3:A:703:936:H291	3:A:703:936:H192	1.75	0.41
1:B:29:LEU:CD1	1:B:211:LEU:HD13	2.50	0.41
1:B:221:TYR:HB2	1:B:273:LYS:HA	2.01	0.41
1:C:233:GLN:HA	1:C:233:GLN:NE2	2.33	0.41
1:B:272:ARG:O	1:B:273:LYS:HB2	2.19	0.41
1:C:141:LYS:NZ	1:C:215:TRP:O	2.51	0.41
1:D:229:ASP:OD1	1:D:239:LYS:HG2	2.20	0.41
1:D:54:ASP:OD2	1:D:118:ARG:HD2	2.20	0.41
1:C:143:VAL:HG22	1:C:184:LEU:HD12	2.01	0.41
1:C:138:ARG:HA	1:C:180:GLY:O	2.19	0.41
1:A:54:ASP:OD1	1:A:118:ARG:NH1	2.47	0.41
1:C:255:VAL:H	1:C:267:ASN:ND2	2.19	0.41
1:B:31:LYS:HZ1	1:B:208:LYS:NZ	2.19	0.41
1:A:108:VAL:HG13	1:A:109:PRO:HD2	2.03	0.41
1:B:154:MET:CE	1:B:161:HIS:NE2	2.83	0.41
1:B:245:GLU:O	1:B:249:LYS:HE3	2.20	0.41
1:A:92:GLU:HG3	1:A:124:PHE:HE1	1.86	0.41
1:A:167:ILE:O	1:A:170:PRO:HD2	2.21	0.41
1:A:112:LEU:HD22	1:A:116:PHE:CE2	2.56	0.41
1:C:101:PHE:CZ	1:C:108:VAL:HG12	2.56	0.41
1:B:104:GLN:HA	2:B:602:FAD:N5	2.35	0.40
1:B:236:PHE:O	1:D:160:ILE:HG13	2.22	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:133:ASP:CG	1:C:225:SER:H	2.25	0.40
1:D:225:SER:HB2	1:D:230:LEU:HD11	2.02	0.40
1:B:154:MET:CE	1:B:160:ILE:HD11	2.49	0.40
1:A:214:ILE:HG23	1:A:215:TRP:N	2.36	0.40
1:B:101:PHE:CZ	1:B:146:ILE:HG12	2.56	0.40
1:B:7:ILE:HG21	1:B:22:LYS:HG2	2.02	0.40
1:C:54:ASP:OD2	1:C:118:ARG:HD2	2.21	0.40
1:C:128:TYR:HB3	3:C:701:936:N34	2.36	0.40
1:D:68:PRO:O	1:D:72:VAL:HG23	2.22	0.40
1:C:11:HIS:CE1	1:C:16:SER:HB3	2.56	0.40
1:C:211:LEU:HA	1:C:214:ILE:HB	2.02	0.40
1:D:240:LYS:HD2	1:D:240:LYS:N	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	270/273 (99%)	255 (94%)	14 (5%)	1 (0%)	43	25
1	B	269/273 (98%)	256 (95%)	12 (4%)	1 (0%)	43	25
1	C	269/273 (98%)	255 (95%)	13 (5%)	1 (0%)	43	25
1	D	270/273 (99%)	255 (94%)	13 (5%)	2 (1%)	30	13
All	All	1078/1092 (99%)	1021 (95%)	52 (5%)	5 (0%)	38	19

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3	ARG
1	C	272	ARG
1	B	272	ARG
1	D	230	LEU
1	D	191	SER

5.3.2 Protein sidechains ⓘ

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	226/227 (100%)	215 (95%)	11 (5%)	35	15
1	B	226/227 (100%)	216 (96%)	10 (4%)	39	18
1	C	226/227 (100%)	220 (97%)	6 (3%)	57	39
1	D	226/227 (100%)	220 (97%)	6 (3%)	57	39
All	All	904/908 (100%)	871 (96%)	33 (4%)	45	26

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	29	LEU
1	A	47	ASN
1	A	48	PRO
1	A	50	ILE
1	A	112	LEU
1	A	158	GLN
1	A	160	ILE
1	A	202	GLN
1	A	241	GLU
1	A	272	ARG
1	B	14	ARG
1	B	22	LYS
1	B	47	ASN
1	B	50	ILE
1	B	112	LEU
1	B	158	GLN
1	B	200	ARG
1	B	240	LYS
1	B	241	GLU
1	B	273	LYS
1	C	47	ASN
1	C	48	PRO
1	C	158	GLN
1	C	209	LYS
1	C	233	GLN

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Mol	Chain	Res	Type
1	C	273	LYS
1	D	14	ARG
1	D	52	ARG
1	D	158	GLN
1	D	189	THR
1	D	202	GLN
1	D	240	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	ASN
1	A	79	HIS
1	A	158	GLN
1	A	172	GLN
1	A	194	HIS
1	A	202	GLN
1	A	233	GLN
1	A	267	ASN
1	A	268	GLN
1	B	47	ASN
1	B	64	ASN
1	B	79	HIS
1	B	172	GLN
1	B	194	HIS
1	B	202	GLN
1	B	231	ASN
1	B	267	ASN
1	B	268	GLN
1	C	47	ASN
1	C	64	ASN
1	C	66	GLN
1	C	158	GLN
1	C	172	GLN
1	C	233	GLN
1	C	267	ASN
1	C	268	GLN
1	D	47	ASN
1	D	66	GLN
1	D	79	HIS
1	D	158	GLN
1	D	172	GLN

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Mol	Chain	Res	Type
1	D	233	GLN
1	D	243	GLN
1	D	267	ASN
1	D	268	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	FAD	A	601	-	58,58,58	2.41	18 (31%)	85,89,89	1.39	11 (12%)
3	936	A	703	-	28,28,28	4.27	11 (39%)	39,41,41	2.20	12 (30%)
2	FAD	B	602	-	58,58,58	2.37	18 (31%)	85,89,89	1.37	11 (12%)
3	936	B	704	-	28,28,28	4.27	10 (35%)	39,41,41	2.26	11 (28%)
2	FAD	C	603	-	58,58,58	2.43	18 (31%)	85,89,89	1.39	11 (12%)
3	936	C	701	-	28,28,28	4.41	10 (35%)	39,41,41	2.22	10 (25%)
2	FAD	D	604	-	58,58,58	2.39	19 (32%)	85,89,89	1.37	10 (11%)
3	936	D	702	-	28,28,28	4.44	8 (28%)	39,41,41	2.28	11 (28%)

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	A	601	-	-	0/34/50/50	0/1/6/6
3	936	A	703	-	-	0/9/27/27	0/1/3/3
2	FAD	B	602	-	-	0/34/50/50	0/1/6/6
3	936	B	704	-	-	0/9/27/27	0/1/3/3
2	FAD	C	603	-	-	0/34/50/50	0/1/6/6
3	936	C	701	-	-	0/9/27/27	0/1/3/3
2	FAD	D	604	-	-	0/34/50/50	0/1/6/6
3	936	D	702	-	-	0/9/27/27	0/1/3/3

All (112) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	701	936	C3-C2	-16.45	1.35	1.50
3	D	702	936	C3-C2	-16.04	1.35	1.50
3	A	703	936	C3-C2	-15.34	1.36	1.50
3	B	704	936	C3-C2	-15.19	1.36	1.50
3	D	702	936	O36-N34	11.07	1.41	1.23
3	B	704	936	O36-N34	10.70	1.40	1.23
3	C	701	936	O36-N34	10.33	1.40	1.23
3	A	703	936	O36-N34	9.84	1.39	1.23
2	B	602	FAD	P-O3P	-8.87	1.43	1.59
2	A	601	FAD	P-O3P	-8.76	1.44	1.59
3	D	702	936	O35-N34	8.42	1.41	1.25
2	C	603	FAD	P-O3P	-8.41	1.44	1.59
3	C	701	936	O35-N34	8.10	1.40	1.25
2	D	604	FAD	P-O3P	-8.08	1.45	1.59
3	A	703	936	O35-N34	7.88	1.40	1.25
3	B	704	936	O35-N34	7.79	1.40	1.25
3	A	703	936	C27-N34	-6.96	1.35	1.46
2	A	601	FAD	C9A-N10	6.94	1.49	1.38
3	D	702	936	C27-N34	-6.81	1.35	1.46
2	C	603	FAD	C9A-N10	6.68	1.48	1.38
2	D	604	FAD	C9A-N10	6.58	1.48	1.38
3	C	701	936	C27-N34	-6.36	1.36	1.46
2	B	602	FAD	C9A-N10	6.29	1.48	1.38
3	B	704	936	C27-N34	-6.09	1.37	1.46
2	C	603	FAD	C4X-C10	5.23	1.50	1.40
2	D	604	FAD	C4X-C10	4.86	1.49	1.40
2	C	603	FAD	C4-C4X	4.85	1.49	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	604	FAD	O4B-C1B	4.76	1.48	1.41
2	B	602	FAD	PA-O2A	-4.60	1.34	1.55
2	A	601	FAD	C4X-C10	4.56	1.48	1.40
2	B	602	FAD	C4X-C10	4.51	1.48	1.40
2	A	601	FAD	PA-O2A	-4.41	1.35	1.55
2	D	604	FAD	C1'-C2'	4.41	1.55	1.51
2	C	603	FAD	PA-O2A	-4.38	1.35	1.55
2	A	601	FAD	O4B-C1B	4.36	1.48	1.41
2	D	604	FAD	PA-O2A	-4.32	1.35	1.55
2	C	603	FAD	C2-N3	4.08	1.45	1.37
2	C	603	FAD	O4B-C1B	4.03	1.47	1.41
2	B	602	FAD	O4B-C1B	3.97	1.47	1.41
3	B	704	936	C4-C5	-3.76	1.36	1.46
3	C	701	936	C4-C5	-3.67	1.37	1.46
2	A	601	FAD	C1'-C2'	3.66	1.55	1.51
3	D	702	936	C4-C5	-3.58	1.37	1.46
3	A	703	936	C4-C5	-3.56	1.37	1.46
2	B	602	FAD	P-O2P	-3.55	1.39	1.55
2	D	604	FAD	C2-N3	3.53	1.44	1.37
2	B	602	FAD	C2-N3	3.45	1.44	1.37
2	B	602	FAD	C4-C4X	3.44	1.46	1.41
2	A	601	FAD	P-O2P	-3.39	1.39	1.55
2	A	601	FAD	C4-C4X	3.38	1.46	1.41
2	C	603	FAD	C4A-N3A	3.38	1.40	1.35
2	D	604	FAD	P-O2P	-3.34	1.40	1.55
2	A	601	FAD	C2-N3	3.33	1.43	1.37
2	C	603	FAD	P-O2P	-3.31	1.40	1.55
2	B	602	FAD	C1'-C2'	3.29	1.54	1.51
3	A	703	936	C19-C9	3.29	1.56	1.51
2	D	604	FAD	C4-C4X	3.26	1.46	1.41
2	A	601	FAD	C4A-N3A	3.16	1.40	1.35
2	B	602	FAD	PA-O5B	-3.10	1.45	1.59
2	D	604	FAD	C4A-N3A	2.99	1.40	1.35
2	A	601	FAD	PA-O5B	-2.98	1.45	1.59
3	B	704	936	C19-C9	2.96	1.56	1.51
3	C	701	936	C19-C9	2.93	1.56	1.51
2	D	604	FAD	O5'-C5'	2.83	1.56	1.44
2	C	603	FAD	O5'-C5'	2.79	1.56	1.44
2	B	602	FAD	C6-C5X	2.78	1.45	1.41
2	A	601	FAD	O5'-C5'	2.73	1.56	1.44
2	B	602	FAD	C2A-N1A	2.67	1.39	1.33
2	C	603	FAD	C2A-N1A	2.65	1.39	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	603	FAD	C8-C7	2.63	1.48	1.40
2	B	602	FAD	C4A-N3A	2.61	1.39	1.35
2	C	603	FAD	C5X-N5	2.59	1.39	1.35
2	D	604	FAD	PA-O5B	-2.58	1.47	1.59
2	D	604	FAD	C5X-N5	2.58	1.39	1.35
2	B	602	FAD	C5X-N5	2.58	1.39	1.35
2	C	603	FAD	C1'-C2'	2.55	1.54	1.51
2	B	602	FAD	C8-C7	2.54	1.48	1.40
2	C	603	FAD	PA-O5B	-2.53	1.47	1.59
2	A	601	FAD	C2A-N1A	2.52	1.38	1.33
2	B	602	FAD	O5'-C5'	2.52	1.55	1.44
2	A	601	FAD	C2A-N3A	2.49	1.37	1.32
3	D	702	936	C29-C24	2.46	1.43	1.38
2	D	604	FAD	C2A-N1A	2.45	1.38	1.33
2	B	602	FAD	C2A-N3A	2.44	1.37	1.32
2	A	601	FAD	C8-C7	2.43	1.47	1.40
3	A	703	936	C4-C9	2.43	1.44	1.39
2	D	604	FAD	C8-C7	2.43	1.47	1.40
3	B	704	936	C4-C9	2.41	1.44	1.39
3	B	704	936	O20-C24	2.41	1.43	1.37
2	A	601	FAD	C6-C5X	2.39	1.44	1.41
3	A	703	936	O20-C24	2.38	1.43	1.37
2	D	604	FAD	C2A-N3A	2.36	1.36	1.32
2	C	603	FAD	C4-N3	2.35	1.41	1.37
2	C	603	FAD	O4B-C4B	2.32	1.50	1.45
2	C	603	FAD	C2A-N3A	2.29	1.36	1.32
2	A	601	FAD	C5X-N5	2.27	1.38	1.35
3	A	703	936	C1-C2	2.25	1.50	1.44
2	D	604	FAD	O4B-C4B	2.25	1.50	1.45
3	B	704	936	C29-C24	2.20	1.43	1.38
2	D	604	FAD	C6-C5X	2.17	1.44	1.41
2	B	602	FAD	C4-N3	2.16	1.40	1.37
3	C	701	936	O20-C24	2.15	1.42	1.37
3	D	702	936	C4-C9	2.12	1.44	1.39
3	C	701	936	C29-C24	2.11	1.43	1.38
3	C	701	936	C1-C2	2.10	1.49	1.44
3	D	702	936	O20-C24	2.09	1.42	1.37
3	B	704	936	C28-C27	2.09	1.43	1.38
2	A	601	FAD	C4-N3	2.08	1.40	1.37
3	A	703	936	O44-C6	2.08	1.40	1.36
3	C	701	936	C4-C9	2.08	1.44	1.39
2	D	604	FAD	C4-N3	2.05	1.40	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	703	936	C29-C24	2.02	1.42	1.38

All (87) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	702	936	O44-C6-C5	8.21	118.42	111.13
3	B	704	936	O44-C6-C5	7.48	117.77	111.13
3	C	701	936	O44-C6-C5	7.14	117.47	111.13
3	A	703	936	O44-C6-C5	7.10	117.43	111.13
2	C	603	FAD	C2-N1-C10	6.06	121.09	114.98
2	A	601	FAD	C2-N1-C10	5.69	120.71	114.98
2	B	602	FAD	C2-N1-C10	5.64	120.66	114.98
2	D	604	FAD	C2-N1-C10	5.51	120.53	114.98
3	A	703	936	C12-N7-C3	5.19	130.18	124.45
3	C	701	936	C12-N7-C3	5.08	130.07	124.45
3	D	702	936	C12-N7-C3	5.08	130.06	124.45
3	B	704	936	C12-N7-C3	5.03	130.01	124.45
3	B	704	936	O35-N34-C27	5.03	124.08	114.42
3	C	701	936	O35-N34-C27	4.95	123.93	114.42
3	D	702	936	O35-N34-C27	4.71	123.46	114.42
3	A	703	936	O35-N34-C27	4.54	123.14	114.42
3	C	701	936	O35-N34-O36	-4.43	112.38	121.35
3	B	704	936	O35-N34-O36	-4.40	112.45	121.35
3	A	703	936	O35-N34-O36	-4.22	112.81	121.35
3	D	702	936	O35-N34-O36	-4.01	113.23	121.35
2	A	601	FAD	C4X-C10-N10	-4.00	118.51	120.51
2	B	602	FAD	O4B-C1B-N9A	-3.99	104.73	108.44
2	B	602	FAD	C4X-C10-N10	-3.96	118.53	120.51
2	D	604	FAD	C4X-C10-N10	-3.94	118.54	120.51
2	D	604	FAD	O4B-C1B-N9A	-3.71	104.99	108.44
2	C	603	FAD	O4B-C1B-N9A	-3.71	104.99	108.44
3	A	703	936	O36-N34-C27	3.68	124.05	119.12
2	C	603	FAD	C4X-C10-N10	-3.61	118.71	120.51
2	A	601	FAD	O4B-C1B-N9A	-3.55	105.14	108.44
3	C	701	936	O36-N34-C27	3.41	123.68	119.12
3	B	704	936	C9-C8-N7	3.40	109.94	106.86
3	A	703	936	C9-C8-N7	3.37	109.91	106.86
3	C	701	936	C9-C8-N7	3.32	109.87	106.86
2	D	604	FAD	O5B-PA-O1A	-3.29	96.50	109.37
3	B	704	936	O36-N34-C27	3.25	123.47	119.12
2	C	603	FAD	O5B-PA-O1A	-3.18	96.93	109.37
3	D	702	936	O36-N34-C27	3.12	123.30	119.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	FAD	O5B-PA-O1A	-3.12	97.15	109.37
2	B	602	FAD	O5B-PA-O1A	-3.04	97.47	109.37
3	D	702	936	C9-C8-N7	3.03	109.61	106.86
3	D	702	936	C19-C9-C8	-2.82	122.29	127.25
2	C	603	FAD	P-O3P-PA	2.80	139.91	131.68
3	C	701	936	O20-C19-C9	2.80	114.11	107.73
3	B	704	936	C19-C9-C8	-2.61	122.67	127.25
2	A	601	FAD	C4B-O4B-C1B	-2.59	106.93	109.75
3	C	701	936	C19-C9-C8	-2.59	122.69	127.25
2	A	601	FAD	N3A-C4A-N9A	2.53	130.01	125.43
2	D	604	FAD	N3A-C4A-N9A	2.45	129.85	125.43
2	D	604	FAD	P-O3P-PA	2.44	138.84	131.68
3	D	702	936	C12-N7-C8	-2.41	121.74	124.09
2	C	603	FAD	N3A-C4A-N9A	2.40	129.77	125.43
2	B	602	FAD	P-O3P-PA	2.40	138.71	131.68
2	A	601	FAD	N3A-C2A-N1A	-2.38	126.72	128.71
3	A	703	936	C19-C9-C8	-2.37	123.08	127.25
2	D	604	FAD	C4B-O4B-C1B	-2.34	107.20	109.75
3	D	702	936	C19-O20-C24	-2.33	111.44	117.66
3	A	703	936	C12-N7-C8	-2.33	121.81	124.09
2	C	603	FAD	C4B-O4B-C1B	-2.32	107.22	109.75
2	B	602	FAD	N3A-C4A-N9A	2.32	129.61	125.43
3	D	702	936	O20-C19-C9	2.31	112.99	107.73
3	A	703	936	C3-N7-C8	-2.31	108.01	110.05
2	D	604	FAD	N3A-C2A-N1A	-2.29	126.79	128.71
3	C	701	936	C12-N7-C8	-2.29	121.85	124.09
3	B	704	936	O20-C19-C9	2.29	112.93	107.73
3	B	704	936	C12-N7-C8	-2.24	121.90	124.09
3	A	703	936	O20-C19-C9	2.23	112.80	107.73
3	C	701	936	C3-N7-C8	-2.22	108.08	110.05
2	C	603	FAD	C5A-C4A-N9A	-2.22	103.96	107.16
3	B	704	936	C3-N7-C8	-2.21	108.09	110.05
2	C	603	FAD	N3A-C2A-N1A	-2.21	126.86	128.71
2	A	601	FAD	C5A-C4A-N9A	-2.19	104.00	107.16
2	D	604	FAD	C5A-C4A-N9A	-2.19	104.00	107.16
2	D	604	FAD	C2A-N1A-C6A	2.16	122.67	118.77
2	B	602	FAD	C5A-C4A-N9A	-2.11	104.12	107.16
3	D	702	936	C3-N7-C8	-2.09	108.20	110.05
3	B	704	936	C28-C27-N34	2.08	120.89	118.96
2	C	603	FAD	C2A-N1A-C6A	2.08	122.53	118.77
2	C	603	FAD	C4'-C3'-C2'	-2.08	108.55	113.25
3	A	703	936	C3-C2-C1	2.08	120.28	116.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	FAD	P-O3P-PA	2.05	137.70	131.68
2	B	602	FAD	C4B-O4B-C1B	-2.05	107.53	109.75
3	A	703	936	C1-C6-C5	-2.05	118.65	122.50
2	B	602	FAD	N3A-C2A-N1A	-2.03	127.01	128.71
2	A	601	FAD	C2A-N1A-C6A	2.02	122.42	118.77
2	B	602	FAD	C4'-C3'-C2'	-2.02	108.68	113.25
2	A	601	FAD	C4'-C3'-C2'	-2.02	108.68	113.25
2	B	602	FAD	O2A-PA-O3P	2.02	114.72	105.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.