



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:37 PM GMT

PDB ID : 1GG5  
Title : CRYSTAL STRUCTURE OF A COMPLEX OF HUMAN NAD[P]H-QUINONE OXIDOREDUCTASE AND A CHEMOTHERAPEUTIC DRUG (E09) AT 2.5 Å RESOLUTION  
Authors : Faig, M.; Bianchet, M.A.; Winski, S.; Hargreaves, R.; Moody, C.J.; Hudnott, A.R.; Ross, D.; Amzel, L.M.  
Deposited on : 2000-07-12  
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

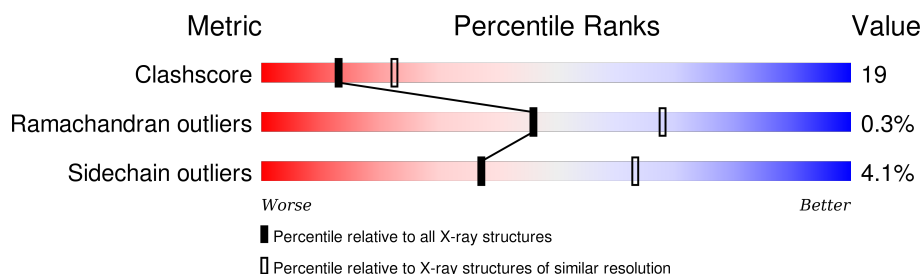
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	273	<div> <div>61%</div> <div>37%</div> <div>.</div> </div>
1	B	273	<div> <div>64%</div> <div>33%</div> <div>.</div> </div>
1	C	273	<div> <div>66%</div> <div>32%</div> <div>.</div> </div>
1	D	273	<div> <div>63%</div> <div>35%</div> <div>.</div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9092 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 NAD(P)H DEHYDROGENASE [QUINONE] 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	0	0
			2173	1412	365	389	7			
1	B	273	Total	C	N	O	S	0	0	0
			2173	1412	365	389	7			
1	C	273	Total	C	N	O	S	0	0	0
			2173	1412	365	389	7			
1	D	273	Total	C	N	O	S	0	0	0
			2173	1412	365	389	7			

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



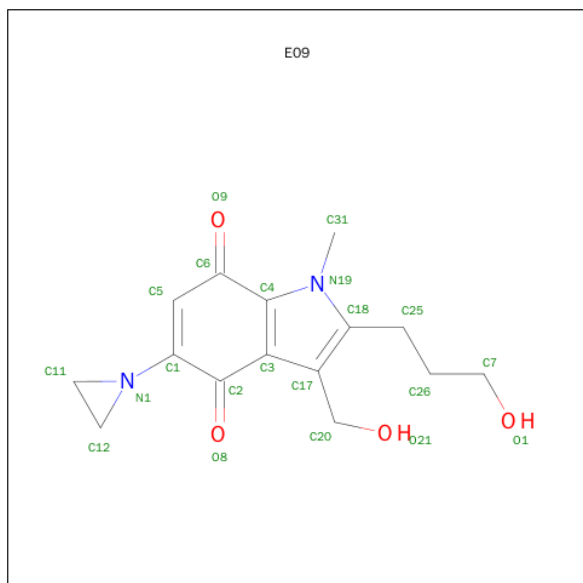
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

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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 3-HYDROXYMETHYL-5-AZIRIDINYL-1METHYL-2-[1H-INDOLE-4,7-DIONE]-PROPANOL (three-letter code: E09) (formula: C<sub>15</sub>H<sub>18</sub>N<sub>2</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	0
			21	15	2	4		
3	D	1	Total	C	N	O	0	0
			21	15	2	4		
3	A	1	Total	C	N	O	0	0
			21	15	2	4		
3	B	1	Total	C	N	O	0	0
			21	15	2	4		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	27	Total	O	0	0
			27	27		
4	B	23	Total	O	0	0
			23	23		
4	C	28	Total	O	0	0
			28	28		

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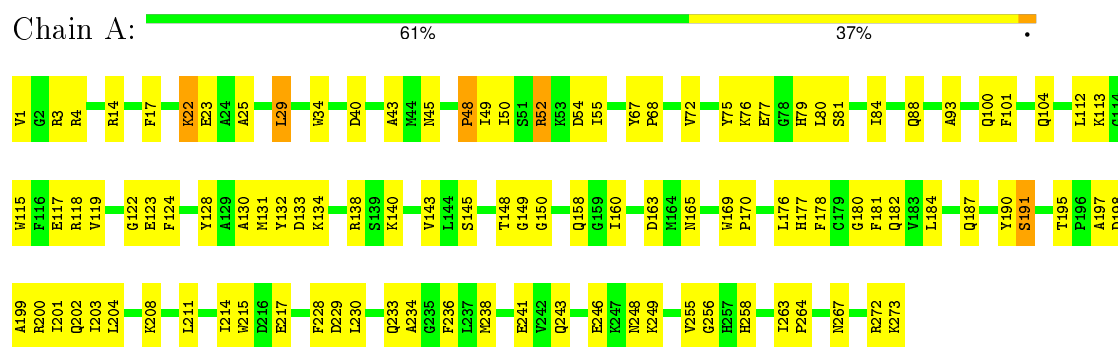
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	26	Total	O	0	0
			26	26		

### 3 Residue-property plots

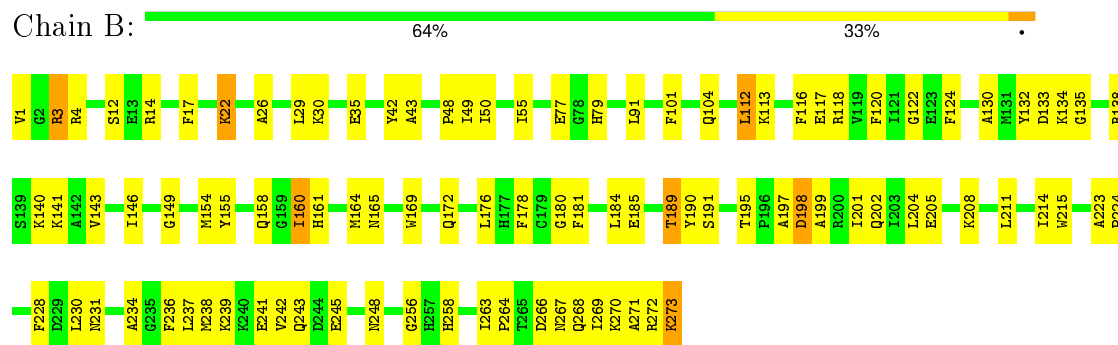
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

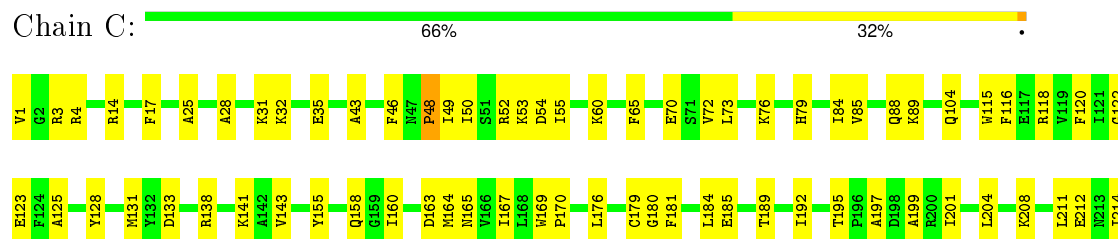
#### • Molecule 1: NAD(P)H DEHYDROGENASE [QUINONE] 1

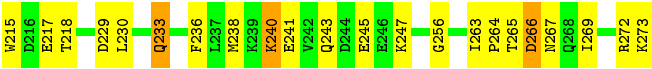


#### • Molecule 1: NAD(P)H DEHYDROGENASE [QUINONE] 1

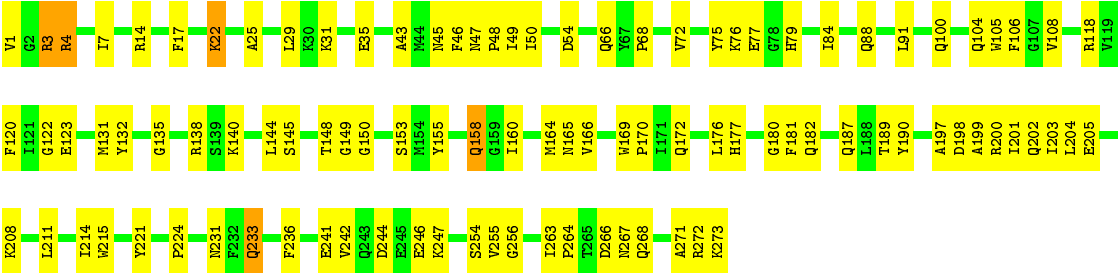


#### • Molecule 1: NAD(P)H DEHYDROGENASE [QUINONE] 1





● Molecule 1: NAD(P)H DEHYDROGENASE [QUINONE] 1



## 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, $\alpha$ , $\beta$ , $\gamma$	56.73 Å   55.45 Å   97.20 Å 76.42°   77.22°   86.45°	Depositor
Resolution (Å)	32.23 – 2.50	Depositor
% Data completeness (in resolution range)	92.3 (32.23-2.50)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
Refinement program	CNS 0.9	Depositor
R, $R_{free}$	0.200 , 0.279	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9092	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP



## 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, E09

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	0.44	0/2231	0.65	0/3012
1	B	0.44	0/2231	0.67	0/3012
1	C	0.44	0/2231	0.65	0/3012
1	D	0.44	0/2231	0.66	1/3012 (0.0%)
All	All	0.44	0/8924	0.65	1/12048 (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	D	182	GLN	N-CA-C	-5.11	97.21	111.00

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	2173	0	2172	103	0
1	B	2173	0	2172	92	0
1	C	2173	0	2172	83	0
1	D	2173	0	2172	94	0
2	A	53	0	31	6	0
2	B	53	0	31	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	53	0	31	4	0
2	D	53	0	31	5	0
3	A	21	0	18	0	0
3	B	21	0	18	6	0
3	C	21	0	18	0	0
3	D	21	0	18	1	0
4	A	27	0	0	1	0
4	B	23	0	0	1	0
4	C	28	0	0	3	0
4	D	26	0	0	1	0
All	All	9092	0	8884	339	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 19.

All (339) 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:PHE:HB2	2:D:604:FAD:H52A	1.39	1.03
1:B:17:PHE:HB2	2:B:602:FAD:H52A	1.52	0.93
1:A:50:ILE:HG22	1:A:118:ARG:HG2	1.50	0.92
1:C:17:PHE:HB2	2:C:603:FAD:H52A	1.50	0.91
1:A:17:PHE:HB2	2:A:601:FAD:H52A	1.51	0.91
1:B:48:PRO:HG3	1:D:49:ILE:HD11	1.56	0.88
1:A:49:ILE:HD11	1:C:48:PRO:HG3	1.59	0.84
1:D:3:ARG:HA	1:D:3:ARG:HE	1.41	0.84
1:A:197:ALA:O	1:A:201:ILE:HG12	1.79	0.83
1:D:1:VAL:HA	1:D:215:TRP:NE1	1.93	0.82
1:D:255:VAL:HG23	1:D:267:ASN:HD22	1.46	0.81
1:D:272:ARG:HB3	1:D:272:ARG:HH11	1.45	0.81
1:B:49:ILE:HD11	1:D:48:PRO:HG3	1.63	0.80
1:C:14:ARG:HH22	1:C:43:ALA:HB2	1.46	0.80
1:C:3:ARG:HA	1:C:3:ARG:HE	1.44	0.80
1:A:160:ILE:HG13	1:C:236:PHE:HB3	1.64	0.79
1:C:240:LYS:HD2	1:C:240:LYS:H	1.46	0.79
1:C:138:ARG:HA	1:C:180:GLY:O	1.83	0.79
1:A:1:VAL:HA	1:A:215:TRP:NE1	1.97	0.78
1:B:3:ARG:HA	1:B:3:ARG:HE	1.47	0.78
1:C:143:VAL:HG22	1:C:184:LEU:HB2	1.68	0.76
1:A:68:PRO:O	1:A:72:VAL:HG23	1.85	0.75
1:B:3:ARG:NE	1:B:3:ARG:HA	2.03	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:204:LEU:O	1:D:208:LYS:HG3	1.87	0.74
1:B:197:ALA:O	1:B:201:ILE:HG12	1.87	0.73
1:D:233:GLN:HE21	1:D:233:GLN:HA	1.55	0.71
1:D:272:ARG:HB3	1:D:272:ARG:NH1	2.05	0.71
1:C:122:GLY:O	1:C:123:GLU:HB2	1.90	0.71
1:D:17:PHE:HB2	2:D:604:FAD:C5B	2.18	0.70
1:B:4:ARG:HG2	1:B:35:GLU:OE1	1.92	0.70
1:A:234:ALA:HB1	4:A:727:HOH:O	1.91	0.70
1:A:3:ARG:NE	1:A:3:ARG:HA	2.07	0.69
1:A:148:THR:HG23	2:A:601:FAD:O2	1.93	0.69
1:B:1:VAL:HA	1:B:215:TRP:NE1	2.09	0.68
1:A:67:TYR:HB3	1:A:68:PRO:HD3	1.76	0.68
1:B:267:ASN:HA	1:B:273:LYS:HE3	1.74	0.68
1:A:48:PRO:HG3	1:C:49:ILE:HD11	1.77	0.67
1:A:3:ARG:HE	1:A:3:ARG:HA	1.58	0.66
1:A:143:VAL:HG22	1:A:184:LEU:HB2	1.77	0.66
1:B:236:PHE:O	1:D:160:ILE:HG13	1.96	0.66
1:C:17:PHE:HB2	2:C:603:FAD:C5B	2.23	0.66
1:A:1:VAL:HA	1:A:215:TRP:HE1	1.60	0.66
1:B:17:PHE:HB2	2:B:602:FAD:C5B	2.25	0.66
1:B:230:LEU:HD23	1:D:160:ILE:CD1	2.26	0.65
1:B:50:ILE:HG22	1:B:118:ARG:HG2	1.78	0.64
1:B:165:ASN:HD21	1:B:266:ASP:HA	1.62	0.64
1:C:50:ILE:HG22	1:C:118:ARG:HG2	1.80	0.64
1:D:201:ILE:O	1:D:205:GLU:HG2	1.98	0.64
1:D:169:TRP:CZ2	1:D:256:GLY:HA3	2.32	0.63
1:A:50:ILE:CG2	1:A:118:ARG:HG2	2.27	0.63
1:A:160:ILE:CG1	1:C:236:PHE:HB3	2.28	0.63
1:B:201:ILE:O	1:B:205:GLU:HG2	1.97	0.63
1:A:204:LEU:O	1:A:208:LYS:HG3	1.98	0.63
1:D:76:LYS:HE3	1:D:123:GLU:HG3	1.79	0.63
1:A:198:ASP:O	1:A:202:GLN:HG2	1.99	0.63
1:C:264:PRO:HB2	1:C:273:LYS:HD2	1.80	0.63
1:B:263:ILE:HD11	1:D:263:ILE:HD11	1.80	0.63
1:C:84:ILE:O	1:C:88:GLN:HG3	1.99	0.63
1:C:204:LEU:O	1:C:208:LYS:HG3	1.98	0.62
1:D:169:TRP:HB3	1:D:170:PRO:HD3	1.82	0.62
1:C:3:ARG:HA	1:C:3:ARG:NE	2.14	0.62
1:B:185:GLU:OE2	1:B:270:LYS:HA	2.01	0.61
1:D:25:ALA:O	1:D:29:LEU:HD23	2.00	0.61
1:C:238:MET:HE2	1:C:243:GLN:HG2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:PHE:HB2	2:A:601:FAD:C5B	2.28	0.60
1:D:267:ASN:HA	1:D:273:LYS:HZ1	1.66	0.60
1:D:1:VAL:HA	1:D:215:TRP:HE1	1.66	0.60
1:B:272:ARG:HH11	1:B:272:ARG:HB3	1.67	0.59
1:D:91:LEU:HD11	1:D:120:PHE:HE1	1.67	0.59
1:A:176:LEU:O	1:A:181:PHE:HB2	2.03	0.59
1:D:264:PRO:HB3	1:D:273:LYS:HD2	1.85	0.59
1:A:40:ASP:HB3	1:A:43:ALA:HB3	1.84	0.59
1:A:160:ILE:HD11	1:C:236:PHE:HD2	1.67	0.58
1:A:202:GLN:HA	1:A:202:GLN:HE21	1.68	0.58
1:A:214:ILE:HD12	1:A:217:GLU:OE2	2.03	0.58
1:B:185:GLU:OE2	1:B:271:ALA:N	2.31	0.58
1:B:189:THR:HG23	4:B:708:HOH:O	2.04	0.58
1:D:76:LYS:CE	1:D:123:GLU:HG3	2.34	0.58
1:A:14:ARG:N	1:A:14:ARG:HD2	2.19	0.58
1:C:14:ARG:HH22	1:C:43:ALA:CB	2.16	0.58
1:B:176:LEU:O	1:B:181:PHE:HB2	2.04	0.58
1:A:258:HIS:HB3	1:A:263:ILE:HG12	1.86	0.57
1:A:54:ASP:OD1	1:A:118:ARG:NH1	2.37	0.57
1:C:76:LYS:HE2	1:C:123:GLU:OE2	2.03	0.57
1:A:72:VAL:HG12	1:A:76:LYS:HD2	1.86	0.57
1:B:264:PRO:HB3	1:B:273:LYS:HD2	1.86	0.56
1:B:202:GLN:HE21	1:B:202:GLN:HA	1.70	0.56
1:B:238:MET:HE3	1:B:243:GLN:HA	1.87	0.56
1:B:236:PHE:HD2	1:D:160:ILE:HD12	1.70	0.56
1:D:155:TYR:HB3	1:D:164:MET:HB2	1.88	0.56
1:A:25:ALA:O	1:A:29:LEU:HD22	2.04	0.56
1:C:169:TRP:HB3	1:C:170:PRO:HD3	1.87	0.56
1:B:113:LYS:O	1:B:117:GLU:HG3	2.06	0.56
1:B:143:VAL:HG22	1:B:184:LEU:HD12	1.87	0.55
1:C:176:LEU:O	1:C:181:PHE:HB2	2.06	0.55
1:A:272:ARG:HG2	1:A:272:ARG:HH11	1.70	0.55
1:A:264:PRO:HB2	1:A:273:LYS:HD2	1.89	0.55
1:B:198:ASP:O	1:B:202:GLN:HG2	2.06	0.55
1:B:143:VAL:CG2	1:B:184:LEU:HD12	2.36	0.55
1:A:75:TYR:CE1	1:A:124:PHE:HB2	2.41	0.55
1:C:54:ASP:OD1	1:C:118:ARG:NH1	2.38	0.55
1:A:160:ILE:CD1	1:C:236:PHE:HB3	2.37	0.55
1:A:200:ARG:NH1	2:A:601:FAD:H1B	2.22	0.55
1:B:160:ILE:HG13	1:D:236:PHE:HB3	1.88	0.55
1:A:4:ARG:NH1	1:A:93:ALA:O	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:167:ILE:O	1:C:170:PRO:HD2	2.06	0.54
1:C:169:TRP:CZ2	1:C:256:GLY:HA3	2.42	0.54
1:D:131:MET:HG2	1:D:132:TYR:CD2	2.42	0.54
1:A:50:ILE:HG13	1:A:50:ILE:O	2.07	0.54
1:B:165:ASN:ND2	1:B:266:ASP:HA	2.22	0.54
1:D:4:ARG:HG2	1:D:35:GLU:HB2	1.90	0.54
1:C:185:GLU:HG3	4:C:713:HOH:O	2.07	0.54
1:A:163:ASP:OD2	1:A:165:ASN:HB2	2.08	0.54
1:B:230:LEU:HD23	1:D:160:ILE:HD13	1.87	0.54
1:B:77:GLU:OE1	1:B:79:HIS:HE1	1.90	0.54
1:A:195:THR:HG22	1:A:199:ALA:HB3	1.89	0.54
1:B:149:GLY:HA3	3:D:702:E09:H252	1.90	0.53
1:C:240:LYS:CD	1:C:240:LYS:H	2.20	0.53
1:B:228:PHE:CG	1:D:160:ILE:HG12	2.43	0.53
1:D:132:TYR:O	1:D:180:GLY:HA2	2.08	0.53
1:B:130:ALA:HB1	1:B:134:LYS:O	2.09	0.53
1:C:197:ALA:O	1:C:201:ILE:HG13	2.09	0.53
1:B:228:PHE:CD2	1:D:160:ILE:HG12	2.44	0.52
1:D:68:PRO:O	1:D:72:VAL:HG23	2.09	0.52
1:D:84:ILE:O	1:D:88:GLN:HG3	2.09	0.52
1:D:104:GLN:HA	2:D:604:FAD:N5	2.25	0.52
1:D:241:GLU:CD	1:D:241:GLU:H	2.13	0.52
1:A:228:PHE:CD2	1:C:160:ILE:HG12	2.45	0.52
1:D:255:VAL:H	1:D:267:ASN:ND2	2.09	0.51
1:C:272:ARG:HH11	1:C:272:ARG:HG2	1.75	0.51
1:A:132:TYR:CD1	1:A:178:PHE:HA	2.45	0.51
1:A:148:THR:CG2	1:A:149:GLY:N	2.74	0.51
1:D:254:SER:HB2	1:D:267:ASN:HD21	1.76	0.51
1:D:54:ASP:OD2	1:D:118:ARG:HD2	2.10	0.51
1:B:202:GLN:NE2	1:B:202:GLN:HA	2.26	0.51
1:B:117:GLU:HG2	1:D:105:TRP:HD1	1.74	0.51
1:A:25:ALA:HA	1:A:211:LEU:CD1	2.40	0.51
1:C:76:LYS:HE3	1:C:123:GLU:HG3	1.91	0.51
1:D:197:ALA:O	1:D:201:ILE:HG13	2.10	0.51
1:C:243:GLN:O	1:C:247:LYS:HG3	2.11	0.51
1:A:130:ALA:HB1	1:A:134:LYS:O	2.10	0.51
1:B:204:LEU:O	1:B:208:LYS:HG3	2.11	0.51
1:B:258:HIS:HB3	1:B:263:ILE:HG12	1.93	0.51
1:A:140:LYS:O	1:A:182:GLN:HG3	2.11	0.51
1:B:77:GLU:HB2	1:B:79:HIS:CE1	2.46	0.50
1:A:148:THR:HG22	1:A:150:GLY:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:271:ALA:O	1:D:272:ARG:HB2	2.11	0.50
1:C:32:LYS:HE3	1:C:212:GLU:HB3	1.93	0.50
3:B:704:E09:H252	1:D:149:GLY:HA3	1.94	0.50
1:A:84:ILE:O	1:A:88:GLN:HG3	2.11	0.50
1:A:236:PHE:O	1:C:160:ILE:HG13	2.12	0.50
1:B:195:THR:HG22	1:B:199:ALA:HB3	1.94	0.50
1:A:169:TRP:CZ2	1:A:256:GLY:HA3	2.46	0.50
1:B:155:TYR:HB3	1:B:164:MET:HB2	1.93	0.50
1:B:113:LYS:NZ	1:D:108:VAL:O	2.44	0.50
1:D:145:SER:HA	1:D:187:GLN:HB3	1.93	0.50
1:B:160:ILE:HD12	1:B:161:HIS:H	1.77	0.50
1:A:255:VAL:HG23	1:A:267:ASN:HD22	1.77	0.49
1:B:267:ASN:HA	1:B:273:LYS:CE	2.42	0.49
1:C:125:ALA:O	1:C:179:CYS:HB3	2.12	0.49
1:A:104:GLN:HA	2:A:601:FAD:N5	2.28	0.49
1:C:128:TYR:N	1:C:128:TYR:CD1	2.81	0.49
1:C:143:VAL:CG2	1:C:184:LEU:HD12	2.43	0.49
3:B:704:E09:H111	1:D:106:PHE:HE1	1.76	0.49
1:B:132:TYR:CD1	1:B:178:PHE:HA	2.48	0.49
1:D:122:GLY:O	1:D:123:GLU:HB2	2.11	0.49
1:B:91:LEU:O	1:B:140:LYS:HE3	2.13	0.49
1:B:237:LEU:HD21	1:D:153:SER:HB2	1.94	0.49
1:D:233:GLN:NE2	1:D:233:GLN:HA	2.25	0.49
1:B:236:PHE:HB3	1:D:160:ILE:HD12	1.94	0.49
1:B:230:LEU:CD2	1:D:160:ILE:HD13	2.43	0.49
1:C:3:ARG:CA	1:C:3:ARG:HE	2.21	0.49
1:A:160:ILE:HD11	1:C:236:PHE:CD2	2.48	0.48
1:A:100:GLN:O	1:A:101:PHE:HB3	2.13	0.48
1:C:4:ARG:HD3	1:C:35:GLU:OE1	2.13	0.48
1:B:169:TRP:CD1	1:D:166:VAL:HG11	2.48	0.48
1:C:1:VAL:HA	1:C:215:TRP:NE1	2.26	0.48
1:B:154:MET:O	1:B:161:HIS:HB2	2.13	0.48
1:B:239:LYS:HB2	1:B:242:VAL:HG23	1.94	0.48
1:D:200:ARG:NH1	2:D:604:FAD:H1B	2.28	0.48
1:C:50:ILE:CG2	1:C:118:ARG:HG2	2.43	0.48
1:D:132:TYR:HA	1:D:177:HIS:O	2.14	0.48
3:B:704:E09:C11	1:D:106:PHE:HE1	2.26	0.48
1:D:104:GLN:HA	2:D:604:FAD:C5X	2.43	0.48
1:C:264:PRO:CB	1:C:273:LYS:HD2	2.42	0.48
1:D:31:LYS:HE3	1:D:31:LYS:HB2	1.66	0.48
1:A:264:PRO:CB	1:A:273:LYS:HD2	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:77:GLU:HB2	1:D:79:HIS:CE1	2.49	0.48
1:A:4:ARG:HD2	1:A:93:ALA:O	2.14	0.48
1:C:240:LYS:HD2	1:C:240:LYS:N	2.22	0.48
1:A:190:TYR:O	1:A:191:SER:C	2.51	0.48
1:D:148:THR:OG1	1:D:190:TYR:HA	2.14	0.47
1:A:122:GLY:O	1:A:123:GLU:HB3	2.14	0.47
1:A:199:ALA:O	1:A:203:ILE:HG13	2.14	0.47
1:A:228:PHE:CG	1:C:160:ILE:HG12	2.50	0.47
1:A:267:ASN:HA	1:A:273:LYS:NZ	2.29	0.47
1:C:4:ARG:HG2	1:C:35:GLU:HB2	1.96	0.47
1:B:165:ASN:OD1	1:B:269:ILE:HG13	2.15	0.47
1:A:132:TYR:HA	1:A:177:HIS:O	2.14	0.47
1:B:138:ARG:HA	1:B:180:GLY:O	2.14	0.47
1:C:55:ILE:HD12	1:C:55:ILE:N	2.30	0.47
1:C:165:ASN:HD21	1:C:266:ASP:HA	1.80	0.47
1:D:244:ASP:HA	1:D:247:LYS:HG3	1.97	0.46
1:C:28:ALA:HA	1:C:31:LYS:HE3	1.97	0.46
1:C:195:THR:HG22	1:C:199:ALA:HB3	1.97	0.46
1:C:214:ILE:HD12	1:C:217:GLU:OE2	2.16	0.46
1:A:132:TYR:O	1:A:180:GLY:HA2	2.15	0.46
1:A:104:GLN:HA	2:A:601:FAD:C5X	2.45	0.46
1:D:43:ALA:C	1:D:45:ASN:H	2.19	0.46
1:D:72:VAL:O	1:D:75:TYR:HB3	2.16	0.46
1:B:267:ASN:OD1	1:B:273:LYS:HE2	2.16	0.46
1:C:267:ASN:HA	1:C:273:LYS:HZ1	1.80	0.46
1:C:165:ASN:ND2	1:C:266:ASP:HA	2.31	0.46
1:A:80:LEU:O	1:A:81:SER:C	2.54	0.45
1:A:263:ILE:HD11	1:C:263:ILE:HD11	1.99	0.45
1:B:190:TYR:O	1:B:191:SER:C	2.55	0.45
1:A:72:VAL:O	1:A:76:LYS:HG3	2.17	0.45
1:B:12:SER:HB3	1:B:42:TYR:CD1	2.52	0.45
3:B:704:E09:H111	1:D:106:PHE:CE1	2.52	0.45
1:D:272:ARG:CB	1:D:272:ARG:NH1	2.79	0.45
1:D:46:PHE:CE2	1:D:47:ASN:O	2.70	0.45
1:C:65:PHE:CE2	1:C:70:GLU:HG3	2.52	0.45
1:D:199:ALA:O	1:D:203:ILE:HG13	2.17	0.45
1:C:264:PRO:HB2	1:C:273:LYS:CD	2.47	0.45
1:B:130:ALA:O	1:B:135:GLY:HA2	2.16	0.45
1:A:122:GLY:O	1:A:123:GLU:CB	2.64	0.45
1:C:60:LYS:HE3	1:C:73:LEU:HD22	1.99	0.44
1:B:14:ARG:HH22	1:B:43:ALA:HB2	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:LYS:O	1:A:117:GLU:HG3	2.17	0.44
1:A:241:GLU:CD	1:A:241:GLU:H	2.20	0.44
1:A:238:MET:CE	1:A:243:GLN:HG2	2.47	0.44
1:D:211:LEU:HA	1:D:214:ILE:HB	1.99	0.44
1:B:104:GLN:HA	2:B:602:FAD:C5X	2.47	0.44
1:C:211:LEU:HA	1:C:214:ILE:HB	1.99	0.44
1:D:242:VAL:O	1:D:246:GLU:HB2	2.18	0.44
1:A:128:TYR:N	1:A:128:TYR:CD1	2.86	0.44
1:D:3:ARG:NE	1:D:3:ARG:HA	2.21	0.44
1:D:7:ILE:HG21	1:D:22:LYS:HG2	2.00	0.44
1:A:236:PHE:HB3	1:C:160:ILE:HG13	2.00	0.44
1:C:229:ASP:O	1:C:236:PHE:HA	2.18	0.43
1:B:239:LYS:O	1:B:243:GLN:HG3	2.17	0.43
1:B:141:LYS:HD3	1:B:184:LEU:HD21	2.00	0.43
1:A:138:ARG:HA	1:A:180:GLY:O	2.18	0.43
1:B:169:TRP:CZ2	1:B:256:GLY:HA3	2.52	0.43
1:A:77:GLU:OE1	1:A:79:HIS:HE1	2.01	0.43
1:C:131:MET:HE1	1:C:236:PHE:CD2	2.54	0.43
1:B:22:LYS:HD2	1:B:22:LYS:C	2.39	0.43
1:A:52:ARG:H	1:A:52:ARG:HG2	1.47	0.43
1:A:214:ILE:HG23	1:A:215:TRP:N	2.33	0.43
1:D:172:GLN:HB2	1:D:268:GLN:NE2	2.32	0.43
1:D:72:VAL:CG1	1:D:76:LYS:HE3	2.48	0.43
1:C:163:ASP:OD2	1:C:165:ASN:HB2	2.18	0.43
1:C:53:LYS:NZ	4:C:705:HOH:O	2.34	0.43
1:C:169:TRP:CE2	1:C:256:GLY:HA3	2.53	0.43
1:A:272:ARG:CZ	1:A:272:ARG:HB3	2.49	0.43
1:D:165:ASN:HD21	1:D:266:ASP:HA	1.83	0.43
1:D:135:GLY:O	1:D:138:ARG:HD3	2.19	0.43
1:B:172:GLN:HB2	1:B:268:GLN:NE2	2.33	0.43
1:D:267:ASN:HA	1:D:273:LYS:HE3	2.01	0.43
1:A:230:LEU:HD23	1:C:160:ILE:HD11	2.00	0.43
1:B:238:MET:CE	1:B:243:GLN:HG2	2.48	0.43
1:C:52:ARG:HG2	1:C:52:ARG:H	1.53	0.43
1:C:233:GLN:HA	1:C:233:GLN:HE21	1.84	0.43
1:A:22:LYS:HD2	1:A:23:GLU:OE1	2.19	0.43
1:D:231:ASN:OD1	1:D:233:GLN:HB3	2.19	0.43
1:B:50:ILE:CG2	1:B:118:ARG:HG2	2.45	0.43
1:D:140:LYS:HB2	1:D:181:PHE:CE1	2.53	0.43
3:B:704:E09:C11	1:D:106:PHE:CE1	3.01	0.43
1:A:145:SER:HA	1:A:187:GLN:HB3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:TRP:N	1:A:170:PRO:HD2	2.34	0.42
1:C:25:ALA:HA	1:C:211:LEU:HD13	2.00	0.42
1:D:45:ASN:CG	1:D:45:ASN:O	2.57	0.42
1:B:272:ARG:NH1	1:B:272:ARG:HB3	2.32	0.42
1:A:267:ASN:CB	1:A:273:LYS:HE3	2.49	0.42
1:A:75:TYR:CZ	1:A:124:PHE:HB2	2.54	0.42
1:C:72:VAL:CG1	1:C:76:LYS:HE3	2.48	0.42
1:A:29:LEU:HB3	1:A:34:TRP:HB2	2.02	0.42
1:A:267:ASN:HB2	1:A:273:LYS:HE3	2.00	0.42
1:A:202:GLN:NE2	1:A:202:GLN:HA	2.32	0.42
1:B:223:ALA:HA	1:B:224:PRO:HD3	1.92	0.42
1:A:1:VAL:HA	1:A:215:TRP:CD1	2.54	0.42
1:A:272:ARG:CG	1:A:272:ARG:HH11	2.30	0.42
1:A:229:ASP:O	1:A:236:PHE:HA	2.19	0.42
1:D:202:GLN:NE2	1:D:202:GLN:HA	2.35	0.42
1:B:55:ILE:N	1:B:55:ILE:HD12	2.34	0.42
1:B:29:LEU:CD2	1:B:211:LEU:HD13	2.50	0.42
1:B:264:PRO:CB	1:B:273:LYS:HD2	2.50	0.42
1:B:238:MET:HE2	1:B:243:GLN:HG2	2.02	0.42
1:C:165:ASN:OD1	1:C:269:ILE:HG13	2.20	0.42
1:A:148:THR:HG22	1:A:150:GLY:N	2.34	0.41
3:B:704:E09:H202	1:D:150:GLY:N	2.35	0.41
1:D:100:GLN:OE1	1:D:145:SER:HB3	2.20	0.41
1:C:55:ILE:HG23	1:C:79:HIS:HD2	1.85	0.41
1:B:241:GLU:H	1:B:241:GLU:CD	2.23	0.41
1:C:192:ILE:CG2	2:C:603:FAD:H5'1	2.50	0.41
1:A:160:ILE:HD11	1:C:236:PHE:HB3	2.03	0.41
1:B:214:ILE:HG23	1:B:215:TRP:N	2.35	0.41
1:B:242:VAL:HA	1:B:245:GLU:OE1	2.20	0.41
1:C:104:GLN:HA	2:C:603:FAD:C5X	2.50	0.41
1:A:255:VAL:O	1:A:258:HIS:HD2	2.04	0.41
1:B:116:PHE:O	1:B:120:PHE:HB2	2.21	0.41
1:C:155:TYR:HB3	1:C:164:MET:HB2	2.02	0.41
1:D:47:ASN:HD22	1:D:48:PRO:HD2	1.86	0.41
1:C:208:LYS:O	1:C:212:GLU:HG3	2.21	0.41
1:A:131:MET:HE2	1:A:236:PHE:CE2	2.55	0.41
1:B:112:LEU:HD22	1:B:116:PHE:CE2	2.55	0.41
1:A:115:TRP:O	1:A:119:VAL:HG23	2.20	0.41
1:A:246:GLU:HA	1:A:249:LYS:HG2	2.01	0.41
1:A:131:MET:HB2	1:A:131:MET:HE3	1.99	0.41
1:C:141:LYS:NZ	1:C:215:TRP:O	2.48	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:GLN:HA	1:A:233:GLN:HE21	1.86	0.41
1:D:144:LEU:HD21	1:D:176:LEU:HD11	2.02	0.41
1:C:230:LEU:HD12	4:C:721:HOH:O	2.20	0.41
1:B:104:GLN:HA	2:B:602:FAD:N5	2.35	0.41
1:D:169:TRP:CE2	1:D:256:GLY:HA3	2.55	0.41
1:D:76:LYS:HE2	1:D:123:GLU:OE2	2.21	0.41
1:B:272:ARG:CB	1:B:272:ARG:NH1	2.84	0.41
1:A:273:LYS:HZ3	1:A:273:LYS:HB3	1.86	0.41
1:B:243:GLN:OE1	1:D:158:GLN:NE2	2.53	0.41
1:B:132:TYR:O	1:B:180:GLY:HA2	2.20	0.41
1:A:55:ILE:HG23	1:A:79:HIS:CD2	2.56	0.41
1:B:122:GLY:C	1:B:124:PHE:H	2.24	0.41
1:C:46:PHE:CZ	1:C:115:TRP:HA	2.56	0.41
1:B:231:ASN:OD1	1:B:234:ALA:N	2.54	0.41
1:B:236:PHE:CD2	1:D:160:ILE:HD12	2.51	0.41
1:D:50:ILE:HG22	1:D:118:ARG:HG2	2.03	0.41
1:D:233:GLN:CA	1:D:233:GLN:HE21	2.24	0.40
1:B:271:ALA:O	1:B:272:ARG:HB2	2.21	0.40
1:A:14:ARG:CD	1:A:14:ARG:N	2.82	0.40
1:B:101:PHE:O	1:B:146:ILE:HA	2.21	0.40
1:A:258:HIS:CD2	1:A:258:HIS:H	2.40	0.40
1:D:246:GLU:HG3	4:D:718:HOH:O	2.19	0.40
1:A:55:ILE:HG23	1:A:79:HIS:HD2	1.87	0.40
1:C:85:VAL:O	1:C:89:LYS:HG3	2.21	0.40
1:A:148:THR:HG22	1:A:149:GLY:N	2.36	0.40
1:B:236:PHE:HB3	1:D:160:ILE:CD1	2.52	0.40
1:D:221:TYR:OH	1:D:224:PRO:HD3	2.21	0.40
1:B:26:ALA:O	1:B:30:LYS:HG3	2.21	0.40
1:C:116:PHE:O	1:C:120:PHE:HB2	2.21	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	271/273 (99%)	251 (93%)	19 (7%)	1 (0%)	39	61
1	B	271/273 (99%)	253 (93%)	18 (7%)	0	100	100
1	C	271/273 (99%)	248 (92%)	21 (8%)	2 (1%)	26	46
1	D	271/273 (99%)	251 (93%)	20 (7%)	0	100	100
All	All	1084/1092 (99%)	1003 (92%)	78 (7%)	3 (0%)	46	68

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	133	ASP
1	C	266	ASP
1	A	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%)	217 (96%)	9 (4%)	38	64
1	B	226/227 (100%)	216 (96%)	10 (4%)	35	60
1	C	226/227 (100%)	217 (96%)	9 (4%)	38	64
1	D	226/227 (100%)	217 (96%)	9 (4%)	38	64
All	All	904/908 (100%)	867 (96%)	37 (4%)	37	63

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LYS
1	A	29	LEU
1	A	45	ASN
1	A	48	PRO
1	A	52	ARG
1	A	112	LEU
1	A	133	ASP
1	A	158	GLN

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Mol	Chain	Res	Type
1	A	248	ASN
1	B	3	ARG
1	B	22	LYS
1	B	112	LEU
1	B	133	ASP
1	B	158	GLN
1	B	160	ILE
1	B	189	THR
1	B	198	ASP
1	B	248	ASN
1	B	273	LYS
1	C	48	PRO
1	C	158	GLN
1	C	189	THR
1	C	218	THR
1	C	233	GLN
1	C	240	LYS
1	C	241	GLU
1	C	245	GLU
1	C	265	THR
1	D	3	ARG
1	D	4	ARG
1	D	14	ARG
1	D	22	LYS
1	D	66	GLN
1	D	158	GLN
1	D	189	THR
1	D	198	ASP
1	D	233	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	ASN
1	A	79	HIS
1	A	158	GLN
1	A	182	GLN
1	A	202	GLN
1	A	233	GLN
1	B	47	ASN
1	B	79	HIS
1	B	172	GLN

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Mol	Chain	Res	Type
1	B	202	GLN
1	B	268	GLN
1	C	47	ASN
1	C	79	HIS
1	C	158	GLN
1	C	194	HIS
1	C	233	GLN
1	D	47	ASN
1	D	79	HIS
1	D	172	GLN
1	D	233	GLN
1	D	267	ASN
1	D	268	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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	-	48,58,58	2.31	18 (37%)	54,89,89	2.09	12 (22%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	E09	A	703	-	19,23,23	2.01	6 (31%)	19,34,34	2.19	6 (31%)
2	FAD	B	602	-	48,58,58	2.28	17 (35%)	54,89,89	2.06	14 (25%)
3	E09	B	704	-	19,23,23	1.88	5 (26%)	19,34,34	2.10	8 (42%)
2	FAD	C	603	-	48,58,58	2.28	16 (33%)	54,89,89	2.08	12 (22%)
3	E09	C	701	-	19,23,23	2.05	7 (36%)	19,34,34	2.18	6 (31%)
2	FAD	D	604	-	48,58,58	2.22	15 (31%)	54,89,89	2.03	12 (22%)
3	E09	D	702	-	19,23,23	2.00	7 (36%)	19,34,34	2.14	8 (42%)

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	-	-	0/30/50/50	0/6/6/6
3	E09	A	703	-	-	0/6/28/28	0/2/3/3
2	FAD	B	602	-	-	0/30/50/50	0/6/6/6
3	E09	B	704	-	-	0/6/28/28	0/2/3/3
2	FAD	C	603	-	-	0/30/50/50	0/6/6/6
3	E09	C	701	-	-	0/6/28/28	0/2/3/3
2	FAD	D	604	-	-	0/30/50/50	0/6/6/6
3	E09	D	702	-	-	0/6/28/28	0/2/3/3

All (91) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	604	FAD	PA-O2A	-4.61	1.35	1.54
2	B	602	FAD	PA-O2A	-4.58	1.35	1.54
2	C	603	FAD	PA-O2A	-4.56	1.35	1.54
2	A	601	FAD	PA-O2A	-4.26	1.36	1.54
2	D	604	FAD	P-O2P	-4.00	1.37	1.54
2	C	603	FAD	P-O2P	-3.89	1.38	1.54
2	A	601	FAD	P-O2P	-3.87	1.38	1.54
2	B	602	FAD	P-O2P	-3.87	1.38	1.54
3	A	703	E09	C26-C25	-3.79	1.33	1.52
3	C	701	E09	C26-C25	-3.57	1.34	1.52
3	B	704	E09	C26-C25	-3.54	1.34	1.52
3	D	702	E09	C26-C25	-3.48	1.35	1.52
2	A	601	FAD	C10-N10	-2.80	1.35	1.39
2	A	601	FAD	PA-O5B	-2.41	1.48	1.59
2	B	602	FAD	PA-O5B	-2.22	1.48	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	602	FAD	C10-N10	-2.05	1.36	1.39
2	A	601	FAD	C2A-N3A	2.02	1.35	1.32
2	C	603	FAD	O4B-C4B	2.02	1.49	1.45
2	A	601	FAD	C8-C7	2.02	1.46	1.41
2	C	603	FAD	C8-C7	2.05	1.46	1.41
2	D	604	FAD	C5B-C4B	2.06	1.58	1.51
2	D	604	FAD	C2A-N3A	2.11	1.35	1.32
2	A	601	FAD	C6-C5X	2.13	1.45	1.41
2	D	604	FAD	C2A-N1A	2.13	1.37	1.33
3	C	701	E09	C5-C6	2.18	1.49	1.44
2	D	604	FAD	C4X-N5	2.18	1.36	1.33
2	A	601	FAD	C5B-C4B	2.18	1.58	1.51
2	A	601	FAD	C5X-N5	2.24	1.38	1.35
3	D	702	E09	C5-C6	2.26	1.50	1.44
2	C	603	FAD	C2A-N1A	2.30	1.38	1.33
2	A	601	FAD	C4A-N3A	2.32	1.39	1.35
2	C	603	FAD	C6-C5X	2.32	1.45	1.41
3	A	703	E09	C11-N1	2.34	1.50	1.46
2	A	601	FAD	C2A-N1A	2.34	1.38	1.33
2	B	602	FAD	C5B-C4B	2.37	1.59	1.51
3	C	701	E09	C11-N1	2.39	1.50	1.46
2	B	602	FAD	C6-C5X	2.41	1.45	1.41
2	C	603	FAD	C5B-C4B	2.41	1.59	1.51
2	C	603	FAD	C5X-N5	2.44	1.39	1.35
2	B	602	FAD	C8-C7	2.56	1.47	1.41
2	B	602	FAD	C2A-N1A	2.58	1.38	1.33
3	B	704	E09	C12-N1	2.59	1.50	1.46
2	D	604	FAD	C8-C7	2.61	1.48	1.41
2	B	602	FAD	C5X-N5	2.62	1.39	1.35
2	B	602	FAD	C4-C4X	2.66	1.46	1.41
3	D	702	E09	C11-N1	2.67	1.50	1.46
2	D	604	FAD	C5X-N5	2.69	1.39	1.35
3	D	702	E09	C12-N1	2.69	1.50	1.46
2	B	602	FAD	C4A-N3A	2.72	1.39	1.35
2	C	603	FAD	C4A-N3A	2.74	1.39	1.35
2	D	604	FAD	C4A-N3A	2.77	1.39	1.35
2	B	602	FAD	C4X-N5	2.78	1.37	1.33
3	C	701	E09	C12-N1	2.80	1.50	1.46
2	D	604	FAD	C4-C4X	2.94	1.47	1.41
3	D	702	E09	C3-C4	3.02	1.46	1.41
3	B	704	E09	C1-N1	3.03	1.48	1.38
3	A	703	E09	C12-N1	3.07	1.51	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	FAD	O5'-C5'	3.09	1.57	1.44
2	D	604	FAD	O5'-C5'	3.17	1.57	1.44
3	B	704	E09	C3-C4	3.19	1.46	1.41
2	B	602	FAD	O5'-C5'	3.21	1.57	1.44
2	C	603	FAD	O5'-C5'	3.22	1.58	1.44
2	A	601	FAD	C4-C4X	3.24	1.47	1.41
2	C	603	FAD	C4X-N5	3.27	1.38	1.33
3	C	701	E09	C3-C4	3.29	1.46	1.41
2	A	601	FAD	C4X-N5	3.31	1.38	1.33
2	C	603	FAD	C4-C4X	3.33	1.47	1.41
3	A	703	E09	C3-C4	3.35	1.46	1.41
3	D	702	E09	C1-N1	3.45	1.50	1.38
3	A	703	E09	C1-N1	3.48	1.50	1.38
3	C	701	E09	C1-N1	3.65	1.51	1.38
2	D	604	FAD	C4-N3	3.92	1.40	1.33
2	B	602	FAD	C4-N3	3.97	1.40	1.33
2	A	601	FAD	C4-N3	4.05	1.40	1.33
3	A	703	E09	C3-C17	4.10	1.48	1.39
2	A	601	FAD	C4X-C10	4.13	1.48	1.41
2	B	602	FAD	C4X-C10	4.14	1.48	1.41
3	B	704	E09	C3-C17	4.19	1.48	1.39
2	C	603	FAD	C4-N3	4.19	1.40	1.33
3	D	702	E09	C3-C17	4.23	1.48	1.39
3	C	701	E09	C3-C17	4.42	1.48	1.39
2	D	604	FAD	C4X-C10	4.46	1.49	1.41
2	C	603	FAD	C4X-C10	4.60	1.49	1.41
2	A	601	FAD	O4B-C1B	4.76	1.47	1.41
2	D	604	FAD	O4B-C1B	4.76	1.47	1.41
2	C	603	FAD	O4B-C1B	4.95	1.47	1.41
2	B	602	FAD	O4B-C1B	5.60	1.48	1.41
2	B	602	FAD	C9A-N10	6.69	1.48	1.38
2	C	603	FAD	C9A-N10	6.83	1.48	1.38
2	D	604	FAD	C9A-N10	6.87	1.48	1.38
2	A	601	FAD	C9A-N10	7.73	1.49	1.38

All (78) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	FAD	C4X-C4-N3	-4.58	117.33	123.59
3	C	701	E09	C5-C1-N1	-4.56	117.95	123.47
2	C	603	FAD	C4X-C4-N3	-4.53	117.39	123.59
2	D	604	FAD	C4X-C4-N3	-4.42	117.54	123.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	602	FAD	C4X-C4-N3	-4.37	117.61	123.59
2	A	601	FAD	C4-C4X-C10	-4.24	117.23	119.94
2	D	604	FAD	N3A-C2A-N1A	-4.14	125.73	128.89
3	A	703	E09	C5-C1-N1	-4.14	118.46	123.47
3	D	702	E09	C5-C1-N1	-4.13	118.47	123.47
3	D	702	E09	C12-N1-C11	-4.10	57.87	60.83
3	A	703	E09	C12-N1-C11	-4.07	57.89	60.83
3	C	701	E09	C12-N1-C11	-4.05	57.91	60.83
3	B	704	E09	C5-C1-N1	-4.01	118.61	123.47
3	C	701	E09	C12-N1-C1	-3.93	113.49	121.80
2	C	603	FAD	N3A-C2A-N1A	-3.93	125.89	128.89
3	A	703	E09	C12-N1-C1	-3.92	113.53	121.80
2	A	601	FAD	N3A-C2A-N1A	-3.90	125.91	128.89
3	B	704	E09	C12-N1-C1	-3.86	113.65	121.80
2	B	602	FAD	C4-C4X-C10	-3.86	117.47	119.94
2	C	603	FAD	C4X-C10-N10	-3.83	118.26	120.52
3	D	702	E09	C12-N1-C1	-3.83	113.72	121.80
2	C	603	FAD	C4-C4X-C10	-3.79	117.52	119.94
3	B	704	E09	C12-N1-C11	-3.59	58.24	60.83
2	B	602	FAD	N3A-C2A-N1A	-3.53	126.19	128.89
2	D	604	FAD	C4X-C10-N10	-3.49	118.46	120.52
2	B	602	FAD	O5B-PA-O1A	-3.48	96.12	109.62
2	D	604	FAD	C4-C4X-C10	-3.46	117.72	119.94
2	A	601	FAD	O5B-PA-O1A	-3.35	96.63	109.62
2	D	604	FAD	O5B-PA-O1A	-3.26	96.97	109.62
2	C	603	FAD	O5B-PA-O1A	-3.22	97.11	109.62
2	B	602	FAD	C4X-C10-N10	-3.17	118.65	120.52
2	A	601	FAD	C4X-C10-N10	-2.87	118.83	120.52
2	A	601	FAD	O4B-C1B-N9A	-2.70	102.45	108.10
2	D	604	FAD	O4B-C1B-N9A	-2.40	103.08	108.10
3	B	704	E09	C3-C4-N19	-2.22	106.05	108.24
3	D	702	E09	C4-C3-C2	-2.11	120.70	122.61
2	B	602	FAD	O5B-C5B-C4B	-2.11	101.35	109.12
2	C	603	FAD	O4B-C1B-N9A	-2.10	103.70	108.10
3	C	701	E09	C4-C3-C2	-2.09	120.72	122.61
3	A	703	E09	C3-C4-N19	-2.08	106.19	108.24
2	B	602	FAD	C4B-O4B-C1B	-2.05	107.46	109.72
3	B	704	E09	C4-C3-C2	-2.02	120.78	122.61
2	B	602	FAD	O4B-C1B-N9A	-2.01	103.88	108.10
3	D	702	E09	C11-C12-N1	2.00	61.06	59.59
3	D	702	E09	C12-C11-N1	2.02	61.07	59.59
2	C	603	FAD	C4A-C5A-N7A	2.08	111.39	109.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	602	FAD	C4A-C5A-N7A	2.18	111.48	109.48
2	A	601	FAD	C2A-N1A-C6A	2.19	122.67	118.77
2	D	604	FAD	C4A-C5A-N7A	2.19	111.49	109.48
2	A	601	FAD	C2B-C1B-N9A	2.20	117.65	114.29
2	B	602	FAD	C2A-N1A-C6A	2.22	122.74	118.77
3	D	702	E09	C25-C26-C7	2.23	121.04	112.98
2	D	604	FAD	C2A-N1A-C6A	2.25	122.78	118.77
3	C	701	E09	C12-C11-N1	2.26	61.25	59.59
3	B	704	E09	C25-C26-C7	2.29	121.24	112.98
3	B	704	E09	C12-C11-N1	2.32	61.29	59.59
2	C	603	FAD	C2A-N1A-C6A	2.35	122.96	118.77
3	A	703	E09	C12-C11-N1	2.48	61.41	59.59
3	B	704	E09	C31-N19-C18	2.90	126.93	124.09
3	C	701	E09	C31-N19-C18	2.99	127.03	124.09
3	D	702	E09	C31-N19-C18	3.11	127.14	124.09
3	A	703	E09	C31-N19-C18	3.30	127.33	124.09
2	C	603	FAD	O3P-P-O5'	3.45	112.09	102.94
2	D	604	FAD	P-O3P-PA	3.60	142.83	132.73
2	B	602	FAD	O3P-P-O5'	3.69	112.72	102.94
2	A	601	FAD	P-O3P-PA	3.83	143.49	132.73
2	D	604	FAD	O3P-P-O5'	3.86	113.19	102.94
2	C	603	FAD	P-O3P-PA	4.07	144.17	132.73
2	B	602	FAD	P-O3P-PA	4.18	144.46	132.73
2	A	601	FAD	O3P-P-O5'	4.30	114.35	102.94
2	A	601	FAD	O3P-PA-O5B	4.47	114.79	102.94
2	D	604	FAD	O3P-PA-O5B	4.75	115.53	102.94
2	C	603	FAD	O3P-PA-O5B	5.05	116.34	102.94
2	B	602	FAD	O3P-PA-O5B	5.16	116.63	102.94
2	D	604	FAD	C4-N3-C2	6.93	121.23	115.25
2	A	601	FAD	C4-N3-C2	7.26	121.52	115.25
2	B	602	FAD	C4-N3-C2	7.41	121.65	115.25
2	C	603	FAD	C4-N3-C2	7.59	121.81	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 26 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	FAD	6	0
2	B	602	FAD	4	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	704	E09	6	0
2	C	603	FAD	4	0
2	D	604	FAD	5	0
3	D	702	E09	1	0

## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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

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

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

### 6.3 Carbohydrates [i](#)

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

### 6.4 Ligands [i](#)

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

### 6.5 Other polymers [i](#)

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