Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 09:43 PM GMT

PDB ID	:	1QBG
Title	:	CRYSTAL STRUCTURE OF HUMAN DT-DIAPHORASE (NAD(P)H OXIDOREDUCTASE)
Authors	:	Skelly, J.V.; Sanderson, M.R.; Suter, D.A.; Baumann, U.; Gregory, D.S.; Bennett, M.; Hobbs, S.M.; Neidle, S.
Deposited on	:	1999-04-20
Resolution	:	2.30 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

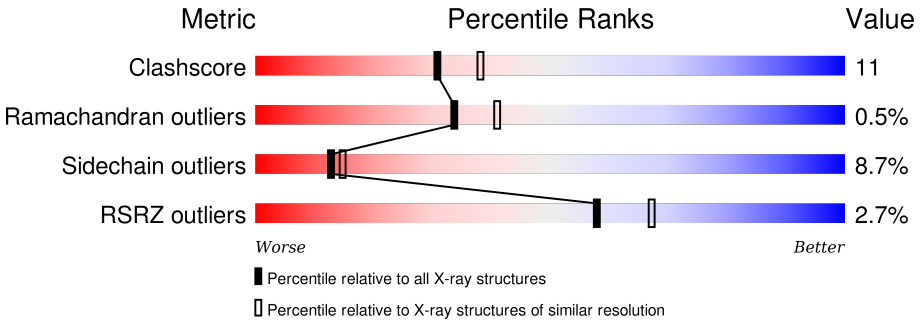
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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. 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 <=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	272	3% 70% 27% ·
1	B	272	3% 68% 29% ·
1	C	272	2% 68% 30% ·
1	D	272	3% 68% 29% ·

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8881 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.

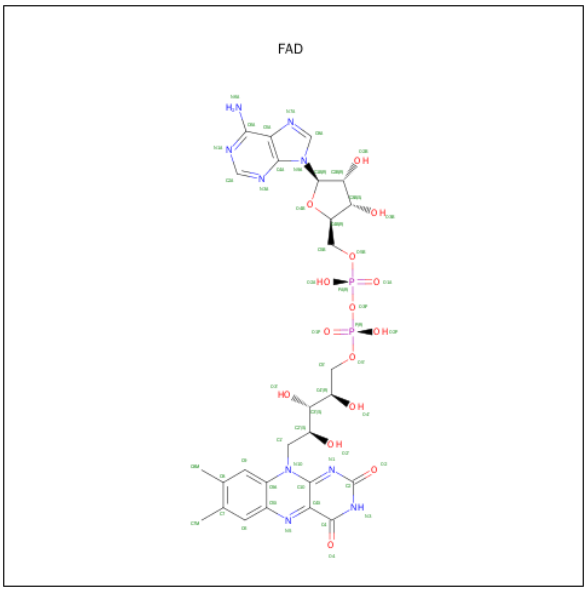
- Molecule 1 is a protein called NAD(P)H DEHYDROGENASE [QUINONE] 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	272	Total	C	N	O	S	0	0	0
			2169	1412	363	387	7			
1	B	271	Total	C	N	O	S	0	0	0
			2162	1407	362	386	7			
1	C	272	Total	C	N	O	S	0	0	0
			2169	1412	363	387	7			
1	D	272	Total	C	N	O	S	0	0	0
			2169	1412	363	387	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	VAL	GLY	conflict	UNP P15559
B	2	VAL	GLY	conflict	UNP P15559
C	2	VAL	GLY	conflict	UNP P15559
D	2	VAL	GLY	conflict	UNP P15559

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



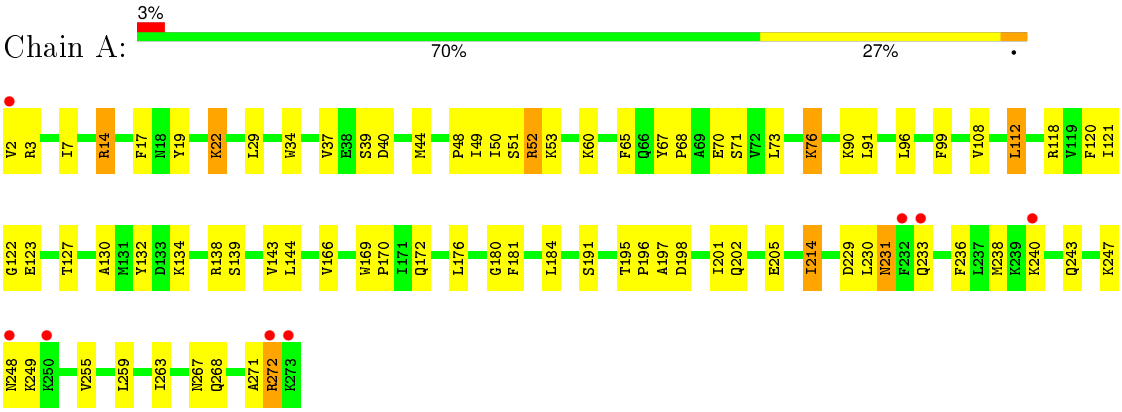
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

3 Residue-property plots (i)

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.

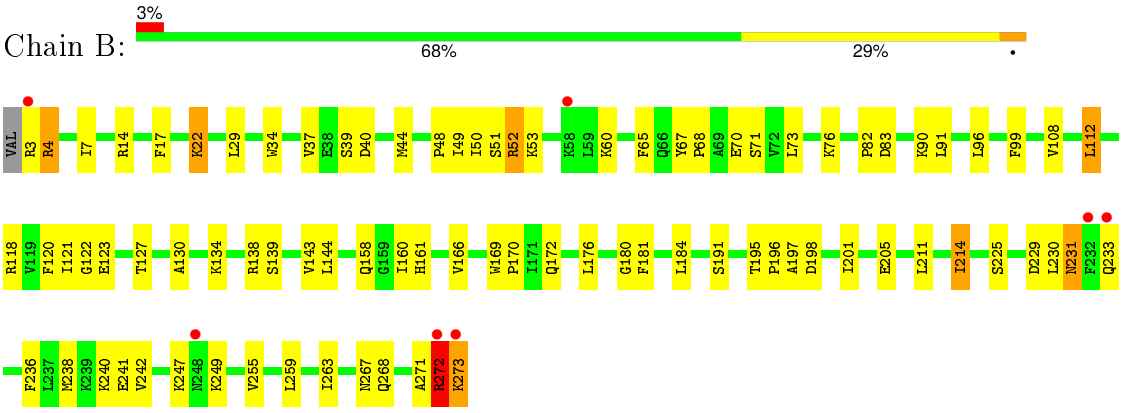
- Molecule 1: NAD(P)H DEHYDROGENASE [QUINONE] 1

Chain A:



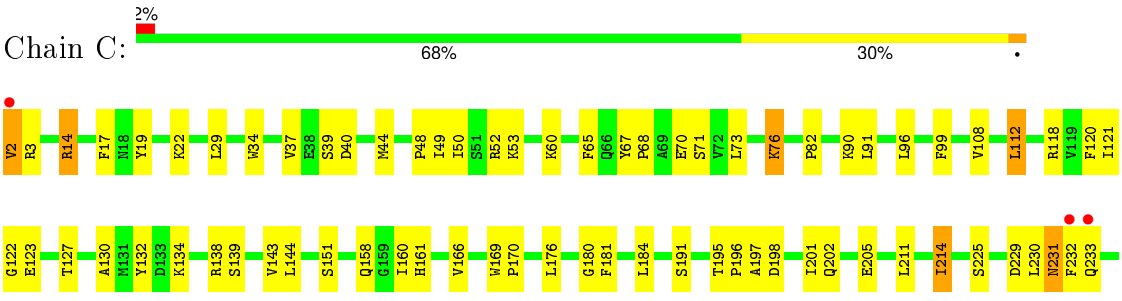
- Molecule 1: NAD(P)H DEHYDROGENASE [QUINONE] 1

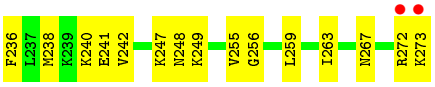
Chain B:



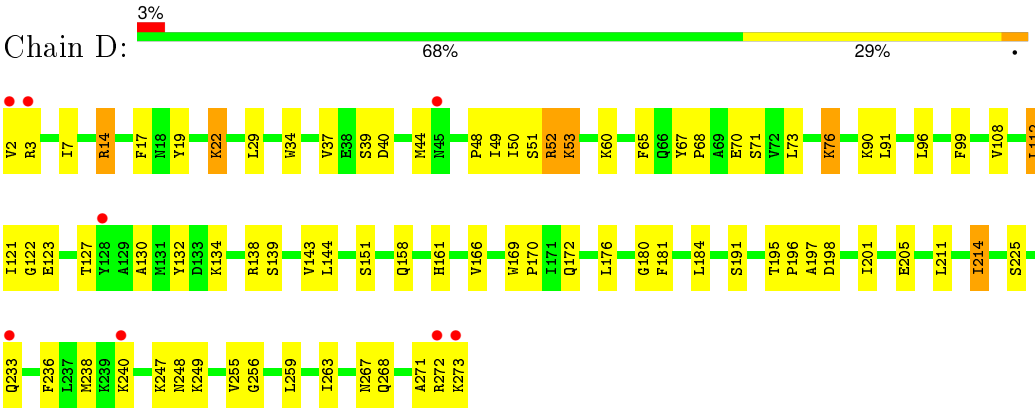
- Molecule 1: NAD(P)H DEHYDROGENASE [QUINONE] 1

Chain C:





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



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	55.88Å 57.49Å 98.77Å 77.10° 76.20° 86.90°	Depositor
Resolution (Å)	40.00 – 2.30 54.27 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (40.00-2.30) 75.8 (54.27-2.20)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.55 (at 2.20Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.238 , 0.280 0.246 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	21.2	Xtriage
Anisotropy	0.135	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 44.6	EDS
Estimated twinning fraction	0.010 for -k,-h,-l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 47496 reflections	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	8881	wwPDB-VP
Average B, all atoms (Å^2)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.42% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.34	0/2227	0.50	0/3009
1	B	0.34	0/2220	0.54	2/2999 (0.1%)
1	C	0.27	0/2227	0.64	3/3009 (0.1%)
1	D	0.28	0/2227	0.47	0/3009
All	All	0.31	0/8901	0.54	5/12026 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	2	VAL	N-CA-C	15.82	153.70	111.00
1	C	2	VAL	CB-CA-C	-12.97	86.75	111.40
1	C	3	ARG	N-CA-CB	-12.80	87.56	110.60
1	B	4	ARG	NE-CZ-NH2	6.57	123.58	120.30
1	B	3	ARG	NE-CZ-NH2	6.40	123.50	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2169	0	2168	54	0
1	B	2162	0	2159	55	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2169	0	2168	51	0
1	D	2169	0	2168	55	0
2	A	53	0	31	2	0
2	B	53	0	31	2	0
2	C	53	0	31	2	0
2	D	53	0	31	2	0
All	All	8881	0	8787	192	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 11.

All (192) 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:C:2:VAL:O	1:C:2:VAL:CG1	1.84	1.24
1:A:243:GLN:OE1	1:B:158:GLN:NE2	1.92	1.03
1:A:243:GLN:OE1	1:B:158:GLN:OE1	1.82	0.96
1:C:2:VAL:HG12	1:C:2:VAL:O	1.12	0.94
1:A:243:GLN:OE1	1:B:158:GLN:CD	2.05	0.94
1:B:50:ILE:HG22	1:B:118:ARG:HG2	1.53	0.89
1:C:50:ILE:HG22	1:C:118:ARG:HG2	1.55	0.89
1:A:50:ILE:HG22	1:A:118:ARG:HG2	1.55	0.88
1:D:50:ILE:HG22	1:D:118:ARG:HG2	1.55	0.88
1:C:49:ILE:HD11	1:D:48:PRO:HG3	1.56	0.87
1:C:48:PRO:HG3	1:D:49:ILE:HD11	1.58	0.86
1:C:255:VAL:HG23	1:C:267:ASN:HD22	1.41	0.85
1:D:255:VAL:HG23	1:D:267:ASN:HD22	1.40	0.85
1:A:48:PRO:HG3	1:B:49:ILE:HD11	1.59	0.85
1:B:255:VAL:HG23	1:B:267:ASN:HD22	1.41	0.85
1:A:49:ILE:HD11	1:B:48:PRO:HG3	1.58	0.84
1:A:255:VAL:HG23	1:A:267:ASN:HD22	1.42	0.81
1:D:271:ALA:O	1:D:272:ARG:HB2	1.88	0.73
1:A:2:VAL:O	1:A:3:ARG:HB2	1.90	0.71
1:A:96:LEU:HD22	1:A:214:ILE:HD13	1.74	0.69
1:B:40:ASP:O	1:B:44:MET:HG3	1.93	0.69
1:D:2:VAL:O	1:D:3:ARG:HB2	1.90	0.69
1:A:166:VAL:HG13	1:B:166:VAL:HG13	1.75	0.69
1:A:76:LYS:HE2	1:A:123:GLU:HG3	1.74	0.68
1:C:76:LYS:HE2	1:C:123:GLU:HG3	1.75	0.68
1:D:40:ASP:O	1:D:44:MET:HG3	1.94	0.67
1:A:40:ASP:O	1:A:44:MET:HG3	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:ASP:HB3	1:D:53:LYS:CE	2.25	0.66
1:B:82:PRO:HG2	1:D:53:LYS:O	1.94	0.66
1:D:76:LYS:HE2	1:D:123:GLU:HG3	1.76	0.66
1:C:166:VAL:HG13	1:D:166:VAL:HG13	1.78	0.65
1:C:40:ASP:O	1:C:44:MET:HG3	1.95	0.65
1:B:134:LYS:HE2	1:B:225:SER:OG	1.97	0.64
1:C:169:TRP:HB3	1:C:170:PRO:HD3	1.81	0.63
1:A:263:ILE:HD11	1:B:263:ILE:HD11	1.81	0.62
1:D:169:TRP:HB3	1:D:170:PRO:HD3	1.80	0.62
1:A:272:ARG:HB2	1:A:272:ARG:HH11	1.64	0.61
1:B:169:TRP:HB3	1:B:170:PRO:HD3	1.83	0.61
1:C:108:VAL:HG13	1:C:112:LEU:HB3	1.83	0.60
1:B:267:ASN:HA	1:B:273:LYS:HE2	1.83	0.60
1:A:191:SER:O	1:A:195:THR:HG23	2.01	0.60
1:B:108:VAL:HG13	1:B:112:LEU:HB3	1.85	0.59
1:B:83:ASP:HB3	1:D:53:LYS:CD	2.32	0.59
1:D:191:SER:O	1:D:195:THR:HG23	2.04	0.58
1:C:191:SER:O	1:C:195:THR:HG23	2.03	0.58
1:D:108:VAL:HG13	1:D:112:LEU:HB3	1.84	0.58
1:A:108:VAL:HG13	1:A:112:LEU:HB3	1.85	0.58
1:C:176:LEU:O	1:C:181:PHE:HB2	2.04	0.58
1:C:122:GLY:O	1:C:123:GLU:HB2	2.04	0.57
1:A:122:GLY:O	1:A:123:GLU:HB2	2.04	0.57
1:B:191:SER:O	1:B:195:THR:HG23	2.04	0.57
1:D:122:GLY:O	1:D:123:GLU:HB2	2.04	0.57
1:D:138:ARG:HA	1:D:180:GLY:O	2.05	0.57
1:D:176:LEU:O	1:D:181:PHE:HB2	2.04	0.56
1:C:138:ARG:HA	1:C:180:GLY:O	2.05	0.56
1:A:176:LEU:O	1:A:181:PHE:HB2	2.04	0.56
1:C:143:VAL:HG22	1:C:184:LEU:HB2	1.87	0.56
1:A:271:ALA:O	1:A:272:ARG:HD3	2.06	0.56
1:A:169:TRP:HB3	1:A:170:PRO:HD3	1.86	0.56
1:A:17:PHE:HB2	2:A:501:FAD:H51A	1.89	0.55
1:B:138:ARG:HA	1:B:180:GLY:O	2.07	0.55
1:B:176:LEU:O	1:B:181:PHE:HB2	2.06	0.55
1:A:138:ARG:HA	1:A:180:GLY:O	2.07	0.54
1:B:96:LEU:HD22	1:B:214:ILE:HD13	1.90	0.54
1:A:143:VAL:HG22	1:A:184:LEU:HB2	1.89	0.54
1:C:96:LEU:HD22	1:C:214:ILE:HD13	1.90	0.53
1:A:271:ALA:O	1:A:272:ARG:HB2	2.08	0.53
1:D:143:VAL:HG22	1:D:184:LEU:HB2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:17:PHE:HB2	2:D:701:FAD:H51A	1.91	0.53
1:B:99:PHE:HB2	1:B:144:LEU:HD23	1.91	0.53
1:D:99:PHE:HB2	1:D:144:LEU:HD23	1.91	0.52
1:C:17:PHE:HB2	2:C:801:FAD:H51A	1.91	0.52
1:A:238:MET:HE3	1:A:259:LEU:HD21	1.91	0.52
1:A:143:VAL:CG2	1:A:184:LEU:HD12	2.40	0.52
1:D:143:VAL:CG2	1:D:184:LEU:HD12	2.40	0.52
1:D:96:LEU:HD22	1:D:214:ILE:HD13	1.92	0.52
1:B:197:ALA:O	1:B:201:ILE:HG12	2.11	0.51
1:A:99:PHE:HB2	1:A:144:LEU:HD23	1.93	0.51
1:D:238:MET:HE1	1:D:259:LEU:HD11	1.93	0.50
1:A:2:VAL:HG12	1:A:3:ARG:H	1.75	0.50
1:C:197:ALA:O	1:C:201:ILE:HG12	2.11	0.50
1:D:2:VAL:HG12	1:D:3:ARG:H	1.76	0.50
1:C:134:LYS:HE2	1:C:225:SER:OG	2.10	0.50
1:A:197:ALA:O	1:A:201:ILE:HG12	2.11	0.50
1:D:197:ALA:O	1:D:201:ILE:HG12	2.12	0.50
1:B:143:VAL:HG22	1:B:184:LEU:HB2	1.94	0.49
1:A:243:GLN:CD	1:B:158:GLN:OE1	2.49	0.49
1:B:17:PHE:HB2	2:B:601:FAD:H51A	1.94	0.49
1:C:143:VAL:CG2	1:C:184:LEU:HD12	2.42	0.49
1:D:172:GLN:HB2	1:D:268:GLN:NE2	2.27	0.48
1:A:132:TYR:OH	1:B:161:HIS:HD2	1.96	0.48
1:B:172:GLN:HB2	1:B:268:GLN:NE2	2.28	0.48
1:C:99:PHE:HB2	1:C:144:LEU:HD23	1.95	0.48
1:B:143:VAL:CG2	1:B:184:LEU:HD12	2.43	0.48
1:B:122:GLY:O	1:B:123:GLU:HB2	2.14	0.48
1:A:29:LEU:O	1:A:34:TRP:HB2	2.14	0.47
1:A:44:MET:HA	1:C:82:PRO:HB2	1.96	0.47
1:D:17:PHE:HB2	2:D:701:FAD:C5B	2.44	0.47
1:B:267:ASN:HA	1:B:273:LYS:CE	2.45	0.47
1:D:29:LEU:O	1:D:34:TRP:HB2	2.15	0.47
1:A:238:MET:HB2	1:B:158:GLN:HB3	1.97	0.46
1:B:29:LEU:O	1:B:34:TRP:HB2	2.15	0.46
1:C:151:SER:HG	1:D:232:PHE:HE1	1.63	0.46
1:D:134:LYS:HE2	1:D:225:SER:OG	2.14	0.46
1:C:29:LEU:O	1:C:34:TRP:HB2	2.15	0.46
1:C:71:SER:HB2	1:C:121:ILE:HD12	1.98	0.46
1:B:83:ASP:HB3	1:D:53:LYS:HE3	1.97	0.46
1:C:263:ILE:HD11	1:D:263:ILE:HD11	1.98	0.45
1:B:17:PHE:HB2	2:B:601:FAD:C5B	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:TYR:N	1:A:68:PRO:HD2	2.32	0.45
1:B:67:TYR:N	1:B:68:PRO:HD2	2.31	0.45
1:C:17:PHE:HB2	2:C:801:FAD:C5B	2.47	0.45
1:A:17:PHE:HB2	2:A:501:FAD:C5B	2.46	0.45
1:D:71:SER:HB2	1:D:121:ILE:HD12	1.99	0.45
1:C:196:PRO:HB2	1:C:198:ASP:OD1	2.17	0.45
1:A:37:VAL:HG11	1:A:90:LYS:HG2	1.99	0.44
1:C:67:TYR:N	1:C:68:PRO:HD2	2.32	0.44
1:C:108:VAL:CG1	1:C:112:LEU:HB3	2.46	0.44
1:D:91:LEU:HD11	1:D:120:PHE:HE1	1.82	0.44
1:D:130:ALA:HB1	1:D:134:LYS:O	2.17	0.44
1:B:108:VAL:CG1	1:B:112:LEU:HB3	2.48	0.44
1:A:65:PHE:CD2	1:A:70:GLU:HG3	2.53	0.44
1:C:91:LEU:HD11	1:C:120:PHE:HE1	1.81	0.44
1:D:67:TYR:N	1:D:68:PRO:HD2	2.32	0.44
1:A:91:LEU:HD11	1:A:120:PHE:HE1	1.83	0.44
1:C:238:MET:HE3	1:C:259:LEU:HD21	2.00	0.43
1:B:91:LEU:HD11	1:B:120:PHE:HE1	1.82	0.43
1:B:267:ASN:HA	1:B:273:LYS:CD	2.47	0.43
1:C:130:ALA:HB1	1:C:134:LYS:O	2.18	0.43
1:C:37:VAL:HG11	1:C:90:LYS:HG2	2.00	0.43
1:D:196:PRO:HB2	1:D:198:ASP:OD1	2.18	0.43
1:A:230:LEU:O	1:A:231:ASN:HB3	2.18	0.43
1:A:272:ARG:HB2	1:A:272:ARG:NH1	2.31	0.43
1:B:196:PRO:HB2	1:B:198:ASP:OD1	2.18	0.43
1:B:52:ARG:H	1:B:52:ARG:HG2	1.41	0.43
1:D:65:PHE:CD2	1:D:70:GLU:HG3	2.53	0.43
1:C:230:LEU:O	1:C:231:ASN:HB3	2.19	0.43
1:B:71:SER:HB2	1:B:121:ILE:HD12	2.00	0.43
1:C:65:PHE:CD2	1:C:70:GLU:HG3	2.54	0.43
1:D:230:LEU:O	1:D:231:ASN:HB3	2.19	0.43
1:B:130:ALA:HB1	1:B:134:LYS:O	2.19	0.43
1:A:229:ASP:O	1:A:236:PHE:HA	2.19	0.43
1:A:52:ARG:H	1:A:52:ARG:HG2	1.42	0.42
1:C:229:ASP:O	1:C:236:PHE:HA	2.19	0.42
1:B:37:VAL:HG11	1:B:90:LYS:HG2	2.00	0.42
1:D:37:VAL:HG11	1:D:90:LYS:HG2	2.01	0.42
1:A:71:SER:HB2	1:A:121:ILE:HD12	2.02	0.42
1:C:169:TRP:CZ2	1:C:256:GLY:HA3	2.54	0.42
1:D:144:LEU:HD12	1:D:172:GLN:HE21	1.85	0.42
1:D:229:ASP:O	1:D:236:PHE:HA	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:ALA:HB1	1:A:134:LYS:O	2.20	0.42
1:B:230:LEU:O	1:B:231:ASN:HB3	2.20	0.42
1:D:211:LEU:HA	1:D:214:ILE:HG12	2.01	0.42
1:B:201:ILE:O	1:B:205:GLU:HG2	2.19	0.42
1:C:132:TYR:OH	1:D:161:HIS:HD2	2.03	0.42
1:B:238:MET:HE1	1:B:259:LEU:HD11	2.01	0.42
1:B:271:ALA:O	1:B:272:ARG:C	2.58	0.42
1:C:238:MET:HE1	1:C:259:LEU:HD11	2.02	0.42
1:D:52:ARG:HG2	1:D:52:ARG:H	1.42	0.42
1:A:196:PRO:HB2	1:A:198:ASP:OD1	2.18	0.42
1:C:211:LEU:HA	1:C:214:ILE:HG12	2.02	0.42
1:D:214:ILE:H	1:D:214:ILE:HG13	1.35	0.42
1:A:172:GLN:HB2	1:A:268:GLN:NE2	2.35	0.42
1:B:123:GLU:OE1	1:B:123:GLU:HA	2.20	0.42
1:B:229:ASP:O	1:B:236:PHE:HA	2.19	0.41
1:C:198:ASP:O	1:C:202:GLN:HG2	2.20	0.41
1:A:7:ILE:HG21	1:A:22:LYS:HG2	2.01	0.41
1:B:211:LEU:HA	1:B:214:ILE:HG12	2.02	0.41
1:A:201:ILE:O	1:A:205:GLU:HG2	2.21	0.41
1:A:198:ASP:O	1:A:202:GLN:HG2	2.20	0.41
1:C:232:PHE:HE1	1:D:151:SER:HG	1.66	0.41
1:C:241:GLU:HG2	1:C:242:VAL:H	1.86	0.41
1:C:2:VAL:O	1:C:34:TRP:CH2	2.74	0.41
1:A:144:LEU:HD12	1:A:172:GLN:HE21	1.85	0.41
1:A:65:PHE:CE2	1:A:70:GLU:HG3	2.55	0.41
1:B:7:ILE:HG21	1:B:22:LYS:HG2	2.03	0.41
1:C:214:ILE:H	1:C:214:ILE:HG13	1.36	0.41
1:D:14:ARG:HG3	1:D:19:TYR:CE1	2.56	0.41
1:B:65:PHE:CD2	1:B:70:GLU:HG3	2.55	0.41
1:D:108:VAL:CG1	1:D:112:LEU:HB3	2.48	0.41
1:C:160:ILE:HG13	1:C:160:ILE:O	2.19	0.41
1:A:214:ILE:HG13	1:A:214:ILE:H	1.52	0.41
1:D:201:ILE:O	1:D:205:GLU:HG2	2.20	0.41
1:B:241:GLU:HG2	1:B:242:VAL:H	1.86	0.41
1:C:14:ARG:HG3	1:C:19:TYR:CE1	2.56	0.41
1:C:161:HIS:HD2	1:D:132:TYR:OH	2.04	0.40
1:C:201:ILE:O	1:C:205:GLU:HG2	2.21	0.40
1:D:7:ILE:HG21	1:D:22:LYS:HG2	2.03	0.40
1:A:14:ARG:HG3	1:A:19:TYR:CE1	2.56	0.40
1:D:169:TRP:CZ2	1:D:256:GLY:HA3	2.56	0.40
1:B:160:ILE:HG13	1:B:160:ILE:O	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:65:PHE:CE2	1:D:70:GLU:HG3	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	270/272 (99%)	251 (93%)	18 (7%)	1 (0%)	39	48
1	B	269/272 (99%)	252 (94%)	15 (6%)	2 (1%)	26	31
1	C	270/272 (99%)	254 (94%)	15 (6%)	1 (0%)	39	48
1	D	270/272 (99%)	252 (93%)	17 (6%)	1 (0%)	39	48
All	All	1079/1088 (99%)	1009 (94%)	65 (6%)	5 (0%)	34	41

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	231	ASN
1	B	231	ASN
1	B	272	ARG
1	C	231	ASN
1	D	231	ASN

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	226/227 (100%)	207 (92%)	19 (8%)	14	16
1	B	225/227 (99%)	205 (91%)	20 (9%)	12	14
1	C	226/227 (100%)	206 (91%)	20 (9%)	12	14
1	D	226/227 (100%)	206 (91%)	20 (9%)	12	14
All	All	903/908 (99%)	824 (91%)	79 (9%)	13	15

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	22	LYS
1	A	39	SER
1	A	51	SER
1	A	52	ARG
1	A	53	LYS
1	A	60	LYS
1	A	73	LEU
1	A	76	LYS
1	A	112	LEU
1	A	127	THR
1	A	139	SER
1	A	214	ILE
1	A	233	GLN
1	A	240	LYS
1	A	247	LYS
1	A	248	ASN
1	A	249	LYS
1	A	272	ARG
1	B	4	ARG
1	B	14	ARG
1	B	22	LYS
1	B	39	SER
1	B	51	SER
1	B	52	ARG
1	B	53	LYS
1	B	60	LYS
1	B	73	LEU
1	B	76	LYS
1	B	112	LEU
1	B	127	THR
1	B	139	SER
1	B	214	ILE

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Mol	Chain	Res	Type
1	B	233	GLN
1	B	240	LYS
1	B	247	LYS
1	B	249	LYS
1	B	272	ARG
1	B	273	LYS
1	C	14	ARG
1	C	22	LYS
1	C	39	SER
1	C	52	ARG
1	C	53	LYS
1	C	60	LYS
1	C	73	LEU
1	C	76	LYS
1	C	112	LEU
1	C	127	THR
1	C	139	SER
1	C	158	GLN
1	C	214	ILE
1	C	233	GLN
1	C	240	LYS
1	C	247	LYS
1	C	248	ASN
1	C	249	LYS
1	C	272	ARG
1	C	273	LYS
1	D	14	ARG
1	D	22	LYS
1	D	39	SER
1	D	51	SER
1	D	52	ARG
1	D	53	LYS
1	D	60	LYS
1	D	73	LEU
1	D	76	LYS
1	D	112	LEU
1	D	127	THR
1	D	139	SER
1	D	158	GLN
1	D	214	ILE
1	D	233	GLN
1	D	240	LYS

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Mol	Chain	Res	Type
1	D	247	LYS
1	D	248	ASN
1	D	249	LYS
1	D	273	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	ASN
1	A	161	HIS
1	A	172	GLN
1	A	202	GLN
1	A	231	ASN
1	A	248	ASN
1	A	267	ASN
1	A	268	GLN
1	B	47	ASN
1	B	161	HIS
1	B	172	GLN
1	B	202	GLN
1	B	231	ASN
1	B	267	ASN
1	B	268	GLN
1	C	47	ASN
1	C	161	HIS
1	C	172	GLN
1	C	202	GLN
1	C	231	ASN
1	C	248	ASN
1	C	267	ASN
1	C	268	GLN
1	D	47	ASN
1	D	161	HIS
1	D	172	GLN
1	D	202	GLN
1	D	231	ASN
1	D	248	ASN
1	D	267	ASN
1	D	268	GLN

5.3.3 RNA

There are no RNA molecules 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

4 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	501	-	48,58,58	2.04	11 (22%)	54,89,89	1.82	12 (22%)
2	FAD	B	601	-	48,58,58	2.05	11 (22%)	54,89,89	1.83	12 (22%)
2	FAD	C	801	-	48,58,58	2.03	10 (20%)	54,89,89	1.84	11 (20%)
2	FAD	D	701	-	48,58,58	2.03	10 (20%)	54,89,89	1.83	12 (22%)

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	501	-	-	0/30/50/50	0/6/6/6
2	FAD	B	601	-	-	0/30/50/50	0/6/6/6
2	FAD	C	801	-	-	0/30/50/50	0/6/6/6
2	FAD	D	701	-	-	0/30/50/50	0/6/6/6

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	FAD	PA-O2A	-4.62	1.35	1.54
2	D	701	FAD	PA-O2A	-4.59	1.35	1.54
2	B	601	FAD	PA-O2A	-4.59	1.35	1.54
2	C	801	FAD	PA-O2A	-4.54	1.35	1.54
2	B	601	FAD	P-O2P	-3.73	1.39	1.54
2	D	701	FAD	P-O2P	-3.69	1.39	1.54
2	A	501	FAD	P-O2P	-3.68	1.39	1.54
2	C	801	FAD	P-O2P	-3.62	1.39	1.54
2	A	501	FAD	C5X-N5	2.03	1.38	1.35
2	B	601	FAD	C5X-N5	2.04	1.38	1.35
2	A	501	FAD	C4A-N3A	2.20	1.38	1.35
2	B	601	FAD	C4A-N3A	2.22	1.38	1.35
2	D	701	FAD	C4A-N3A	2.24	1.38	1.35
2	C	801	FAD	C4A-N3A	2.28	1.39	1.35
2	C	801	FAD	C4-C4X	2.32	1.45	1.41
2	D	701	FAD	C8-C7	2.37	1.47	1.41
2	A	501	FAD	C4-C4X	2.38	1.46	1.41
2	C	801	FAD	C8-C7	2.41	1.47	1.41
2	B	601	FAD	C8-C7	2.44	1.47	1.41
2	B	601	FAD	C4-C4X	2.46	1.46	1.41
2	A	501	FAD	C8-C7	2.51	1.47	1.41
2	D	701	FAD	C4-C4X	2.64	1.46	1.41
2	B	601	FAD	C4-N3	2.86	1.38	1.33
2	A	501	FAD	C4-N3	2.90	1.38	1.33
2	D	701	FAD	C4-N3	2.93	1.38	1.33
2	C	801	FAD	C4-N3	3.09	1.38	1.33
2	C	801	FAD	C4X-C10	3.69	1.47	1.41
2	A	501	FAD	C4X-C10	3.74	1.48	1.41
2	A	501	FAD	O5'-C5'	3.84	1.60	1.44
2	B	601	FAD	O5'-C5'	3.87	1.60	1.44
2	D	701	FAD	O5'-C5'	3.87	1.60	1.44
2	C	801	FAD	O5'-C5'	3.88	1.60	1.44
2	D	701	FAD	C4X-C10	3.88	1.48	1.41
2	B	601	FAD	C4X-C10	3.90	1.48	1.41
2	D	701	FAD	C9A-N10	5.56	1.46	1.38
2	A	501	FAD	C9A-N10	5.59	1.46	1.38
2	B	601	FAD	C9A-N10	5.67	1.46	1.38
2	D	701	FAD	O4B-C1B	5.69	1.48	1.41
2	C	801	FAD	C9A-N10	5.71	1.46	1.38
2	B	601	FAD	O4B-C1B	5.72	1.48	1.41
2	C	801	FAD	O4B-C1B	5.78	1.48	1.41
2	A	501	FAD	O4B-C1B	5.96	1.48	1.41

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	801	FAD	C4X-C4-N3	-4.79	117.04	123.59
2	B	601	FAD	C4X-C4-N3	-4.74	117.11	123.59
2	D	701	FAD	C4X-C4-N3	-4.68	117.19	123.59
2	A	501	FAD	C4X-C4-N3	-4.67	117.21	123.59
2	C	801	FAD	N3A-C2A-N1A	-3.68	126.07	128.89
2	D	701	FAD	N3A-C2A-N1A	-3.68	126.08	128.89
2	B	601	FAD	N3A-C2A-N1A	-3.65	126.10	128.89
2	A	501	FAD	N3A-C2A-N1A	-3.64	126.11	128.89
2	D	701	FAD	O4B-C1B-N9A	-3.24	101.32	108.10
2	A	501	FAD	O4B-C1B-N9A	-3.20	101.40	108.10
2	B	601	FAD	O4B-C1B-N9A	-3.17	101.46	108.10
2	C	801	FAD	O4B-C1B-N9A	-3.14	101.52	108.10
2	A	501	FAD	O5B-PA-O1A	-2.92	98.28	109.62
2	C	801	FAD	O5B-PA-O1A	-2.87	98.47	109.62
2	B	601	FAD	O5B-PA-O1A	-2.87	98.47	109.62
2	D	701	FAD	O5B-PA-O1A	-2.86	98.53	109.62
2	D	701	FAD	C4-C4X-C10	-2.41	118.40	119.94
2	D	701	FAD	C4X-C10-N10	-2.39	119.11	120.52
2	A	501	FAD	C4-C4X-C10	-2.38	118.42	119.94
2	B	601	FAD	C4-C4X-C10	-2.37	118.43	119.94
2	C	801	FAD	C4X-C10-N10	-2.35	119.14	120.52
2	C	801	FAD	C4-C4X-C10	-2.28	118.48	119.94
2	A	501	FAD	C4X-C10-N10	-2.24	119.20	120.52
2	C	801	FAD	C5X-C9A-N10	-2.17	115.97	117.62
2	B	601	FAD	C5X-C9A-N10	-2.13	116.00	117.62
2	B	601	FAD	C4X-C10-N10	-2.13	119.26	120.52
2	A	501	FAD	C5X-C9A-N10	-2.11	116.02	117.62
2	D	701	FAD	C5X-C9A-N10	-2.07	116.04	117.62
2	D	701	FAD	O5'-P-O1P	-2.05	101.65	109.62
2	A	501	FAD	O5'-P-O1P	-2.04	101.71	109.62
2	B	601	FAD	C4A-C5A-N7A	2.02	111.34	109.48
2	C	801	FAD	O2A-PA-O3P	2.09	114.58	105.09
2	D	701	FAD	O2A-PA-O3P	2.11	114.64	105.09
2	A	501	FAD	O2A-PA-O3P	2.11	114.65	105.09
2	B	601	FAD	O2A-PA-O3P	2.13	114.78	105.09
2	C	801	FAD	C2A-N1A-C6A	2.29	122.85	118.77
2	A	501	FAD	C2A-N1A-C6A	2.29	122.86	118.77
2	B	601	FAD	O3P-P-O5'	2.32	109.10	102.94
2	D	701	FAD	C2A-N1A-C6A	2.34	122.95	118.77
2	A	501	FAD	O3P-P-O5'	2.35	109.18	102.94
2	B	601	FAD	C2A-N1A-C6A	2.36	122.98	118.77
2	D	701	FAD	O3P-P-O5'	2.37	109.22	102.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	801	FAD	O3P-P-O5'	2.42	109.35	102.94
2	D	701	FAD	C4-N3-C2	7.37	121.61	115.25
2	A	501	FAD	C4-N3-C2	7.41	121.65	115.25
2	B	601	FAD	C4-N3-C2	7.49	121.72	115.25
2	C	801	FAD	C4-N3-C2	7.54	121.76	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	FAD	2	0
2	B	601	FAD	2	0
2	C	801	FAD	2	0
2	D	701	FAD	2	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

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, 95^{th} 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	272/272 (100%)	0.07	8 (2%)	55	64	3, 14, 40, 62	0
1	B	271/272 (99%)	0.10	7 (2%)	59	68	3, 16, 39, 62	0
1	C	272/272 (100%)	0.05	5 (1%)	71	78	2, 15, 40, 62	0
1	D	272/272 (100%)	0.17	9 (3%)	50	59	3, 17, 41, 61	0
All	All	1087/1088 (99%)	0.10	29 (2%)	58	67	2, 16, 40, 62	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	2	VAL	6.9
1	B	273	LYS	5.2
1	D	2	VAL	5.1
1	A	232	PHE	5.0
1	A	273	LYS	4.8
1	B	232	PHE	4.6
1	D	232	PHE	4.5
1	A	2	VAL	4.3
1	C	273	LYS	4.1
1	C	232	PHE	4.1
1	D	233	GLN	3.5
1	D	272	ARG	3.4
1	D	273	LYS	3.4
1	A	248	ASN	3.4
1	C	272	ARG	3.2
1	A	272	ARG	3.2
1	B	272	ARG	3.0
1	B	58	LYS	2.8
1	A	233	GLN	2.6
1	A	240	LYS	2.6
1	D	45	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	128	TYR	2.5
1	B	233	GLN	2.5
1	B	3	ARG	2.5
1	C	233	GLN	2.5
1	B	248	ASN	2.3
1	D	3	ARG	2.3
1	A	250	LYS	2.3
1	D	240	LYS	2.1

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} 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	LLDF	B-factors(Å^2)	Q<0.9
2	FAD	A	501	53/53	0.95	0.15	0.57	6,15,27,29	0
2	FAD	D	701	53/53	0.93	0.13	0.15	4,18,25,28	0
2	FAD	C	801	53/53	0.94	0.12	-0.04	7,18,22,25	0
2	FAD	B	601	53/53	0.94	0.13	-0.07	9,17,26,31	0

6.5 Other polymers (i)

There are no such residues in this entry.