



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 20, 2016 – 06:58 PM EST

PDB ID : 4GDD
Title : Crystal structure of oxidized Aspergillus fumigatus UDP-galactopyranose mutase complexed with NADH
Authors : Tanner, J.J.; Dhatwalia, R.D.; Singh, H.
Deposited on : 2012-07-31
Resolution : 2.75 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20028442
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)	:	rb-20028442

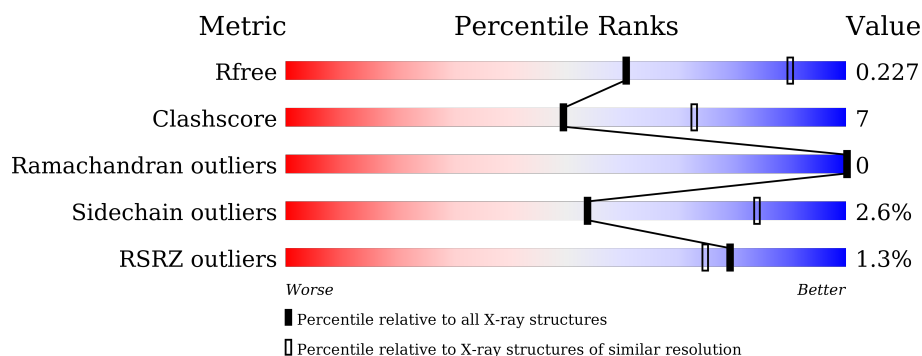
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3340 (2.80-2.72)
Clashscore	102246	3829 (2.80-2.72)
Ramachandran outliers	100387	3767 (2.80-2.72)
Sidechain outliers	100360	3770 (2.80-2.72)
RSRZ outliers	91569	3352 (2.80-2.72)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	513	
1	B	513	
1	C	513	
1	D	513	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SO4	B	605	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 15540 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called UDP-galactopyranose mutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	494	Total	C	N	O	S	0	0	0
			3776	2409	634	712	21			
1	B	494	Total	C	N	O	S	0	0	0
			3777	2406	633	718	20			
1	C	499	Total	C	N	O	S	0	0	0
			3798	2422	641	716	19			
1	D	502	Total	C	N	O	S	0	0	0
			3813	2434	640	719	20			

There are 24 discrepancies between the modelled and reference sequences:

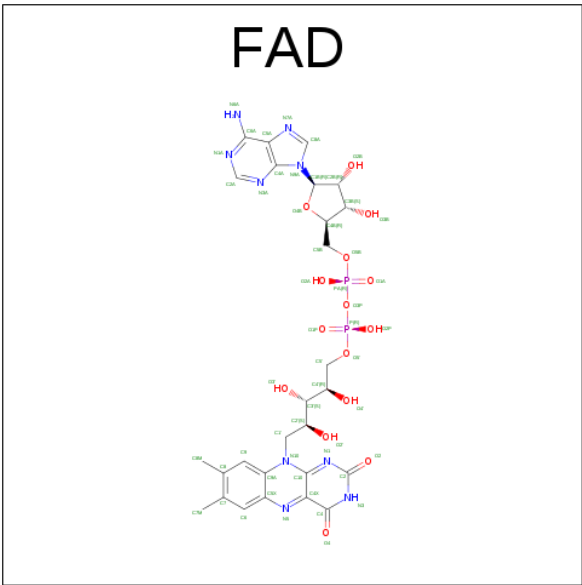
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	ALA	-	EXPRESSION TAG	UNP Q4W1X2
A	-1	ILE	-	EXPRESSION TAG	UNP Q4W1X2
A	0	ALA	-	EXPRESSION TAG	UNP Q4W1X2
A	344	ALA	LYS	ENGINEERED MUTATION	UNP Q4W1X2
A	345	ALA	LYS	ENGINEERED MUTATION	UNP Q4W1X2
A	429	THR	ALA	SEE REMARK 999	UNP Q4W1X2
B	-2	ALA	-	EXPRESSION TAG	UNP Q4W1X2
B	-1	ILE	-	EXPRESSION TAG	UNP Q4W1X2
B	0	ALA	-	EXPRESSION TAG	UNP Q4W1X2
B	344	ALA	LYS	ENGINEERED MUTATION	UNP Q4W1X2
B	345	ALA	LYS	ENGINEERED MUTATION	UNP Q4W1X2
B	429	THR	ALA	SEE REMARK 999	UNP Q4W1X2
C	-2	ALA	-	EXPRESSION TAG	UNP Q4W1X2
C	-1	ILE	-	EXPRESSION TAG	UNP Q4W1X2
C	0	ALA	-	EXPRESSION TAG	UNP Q4W1X2
C	344	ALA	LYS	ENGINEERED MUTATION	UNP Q4W1X2
C	345	ALA	LYS	ENGINEERED MUTATION	UNP Q4W1X2
C	429	THR	ALA	SEE REMARK 999	UNP Q4W1X2
D	-2	ALA	-	EXPRESSION TAG	UNP Q4W1X2
D	-1	ILE	-	EXPRESSION TAG	UNP Q4W1X2
D	0	ALA	-	EXPRESSION TAG	UNP Q4W1X2

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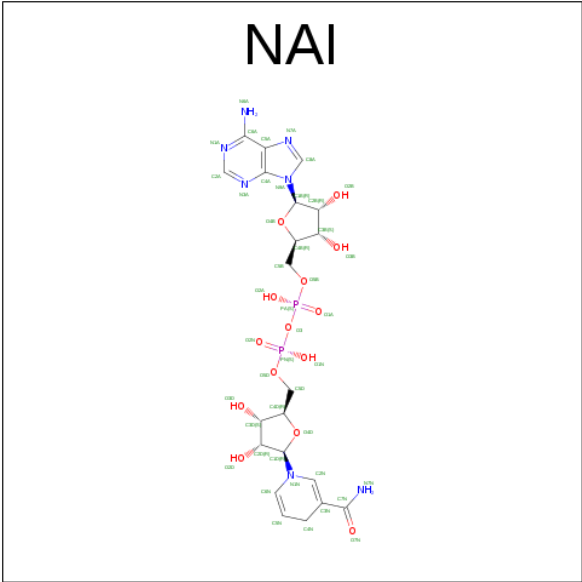
Chain	Residue	Modelled	Actual	Comment	Reference
D	344	ALA	LYS	ENGINEERED MUTATION	UNP Q4W1X2
D	345	ALA	LYS	ENGINEERED MUTATION	UNP Q4W1X2
D	429	THR	ALA	SEE REMARK 999	UNP Q4W1X2

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



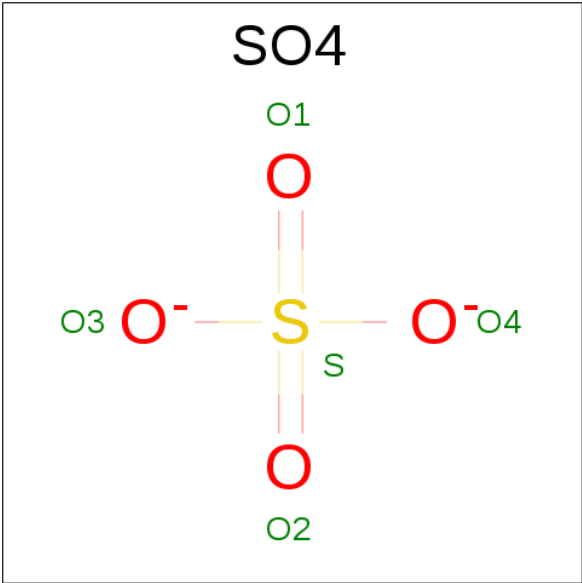
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (three-letter code: NAI) (formula: C₂₁H₂₉N₇O₁₄P₂).



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

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		

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

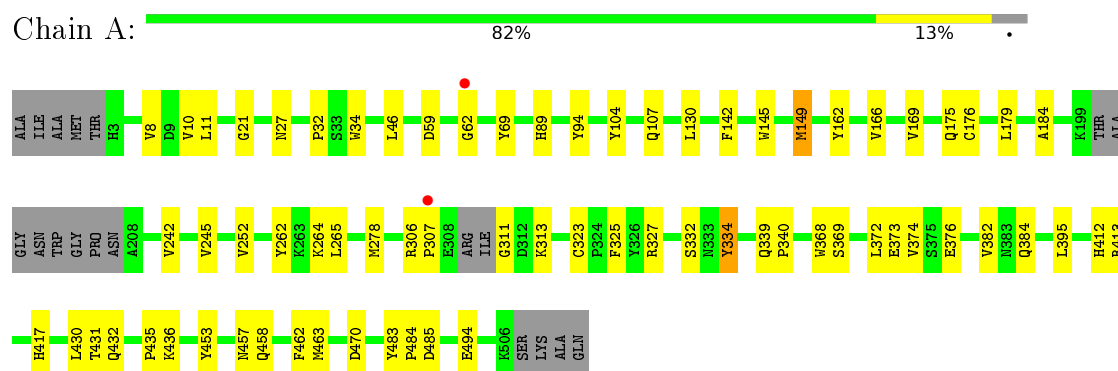
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	4	Total O 4 4	0	0
5	B	7	Total O 7 7	0	0
5	C	7	Total O 7 7	0	0
5	D	8	Total O 8 8	0	0

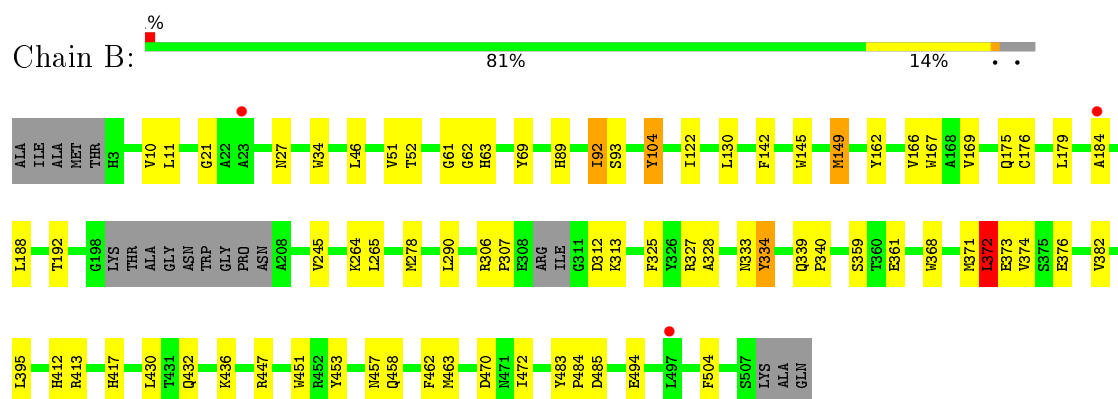
3 Residue-property plots [i](#)

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.

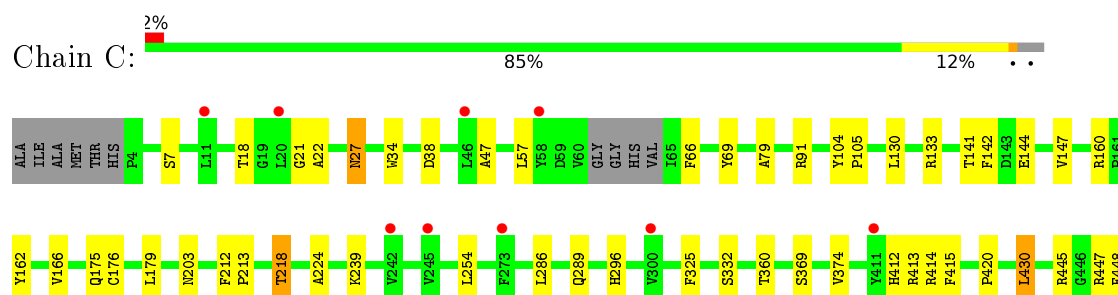
• Molecule 1: UDP-galactopyranose mutase



• Molecule 1: UDP-galactopyranose mutase

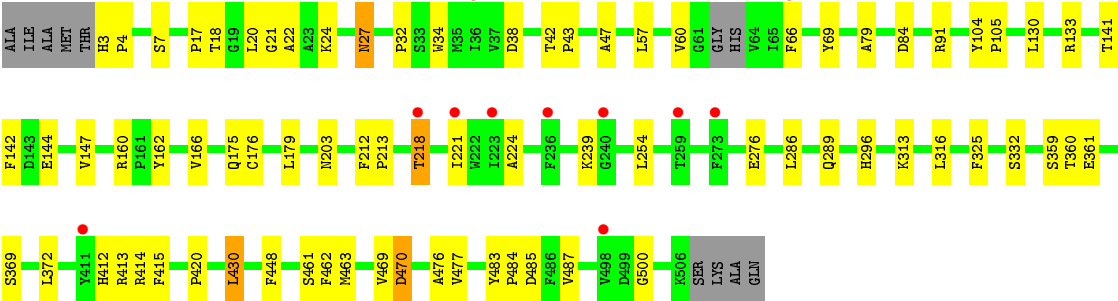
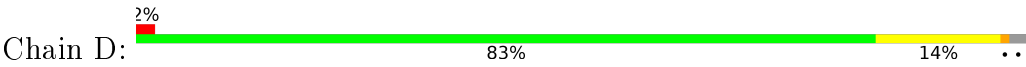


• Molecule 1: UDP-galactopyranose mutase





● Molecule 1: UDP-galactopyranose mutase



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	218.25Å 218.25Å 318.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	162.48 – 2.75 162.48 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.4 (162.48-2.75) 99.3 (162.48-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.73Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.2_432)	Depositor
R, R_{free}	0.212 , 0.245 0.191 , 0.227	Depositor DCC
R_{free} test set	5786 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	51.2	Xtriage
Anisotropy	0.318	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 53.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	15540	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAI, SO4, 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	0.45	0/3869	0.60	0/5276
1	B	0.45	0/3870	0.61	1/5279 (0.0%)
1	C	0.41	0/3894	0.56	0/5314
1	D	0.42	0/3909	0.57	0/5333
All	All	0.43	0/15542	0.59	1/21202 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	372	LEU	CA-CB-CG	6.95	131.28	115.30

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	3776	0	3587	53	0
1	B	3777	0	3583	64	0
1	C	3798	0	3583	46	0
1	D	3813	0	3602	52	0
2	A	53	0	30	3	0
2	B	53	0	30	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	53	0	31	5	0
2	D	53	0	31	3	0
3	A	44	0	27	2	0
3	B	44	0	27	2	0
4	A	10	0	0	0	0
4	B	15	0	0	2	0
4	C	10	0	0	1	0
4	D	15	0	0	0	0
5	A	4	0	0	0	0
5	B	7	0	0	0	0
5	C	7	0	0	0	0
5	D	8	0	0	0	0
All	All	15540	0	14531	214	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 7.

All (214) 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:B:453:TYR:CE1	3:B:602:NAI:H42N	2.04	0.92
1:B:325:PHE:CD1	1:B:372:LEU:HD12	2.07	0.90
1:D:38:ASP:OD1	2:D:601:FAD:H1B	1.83	0.78
1:D:359:SER:OG	1:D:361:GLU:HG2	1.85	0.77
1:A:453:TYR:CE1	3:A:602:NAI:H42N	2.22	0.73
1:D:296:HIS:CE1	1:D:414:ARG:HD2	2.25	0.71
1:B:69:TYR:CD1	1:B:463:MET:HG3	2.26	0.71
1:C:296:HIS:CE1	1:C:414:ARG:HD2	2.25	0.70
1:A:69:TYR:CD1	1:A:463:MET:HG3	2.27	0.70
1:B:62:GLY:HA2	1:B:458:GLN:HE22	1.56	0.70
1:C:38:ASP:OD1	2:C:601:FAD:H1B	1.92	0.70
1:B:327:ARG:NE	1:B:373:GLU:OE2	2.25	0.67
1:A:325:PHE:CD1	1:A:372:LEU:HD22	2.29	0.67
1:A:306:ARG:HD2	1:A:311:GLY:HA2	1.76	0.66
1:B:325:PHE:CG	1:B:372:LEU:HD12	2.31	0.65
1:B:453:TYR:CZ	3:B:602:NAI:H42N	2.31	0.65
1:C:66:PHE:HB2	1:C:91:ARG:NH1	2.11	0.65
1:A:245:VAL:HG22	1:A:252:VAL:HG22	1.78	0.64
1:D:141:THR:OG1	1:D:144:GLU:HG3	1.98	0.64
1:B:122:ILE:HD12	1:B:192:THR:HG22	1.79	0.64
1:C:141:THR:OG1	1:C:144:GLU:HG3	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:GLY:HA2	1:B:458:GLN:NE2	2.12	0.63
1:D:69:TYR:CD1	1:D:463:MET:HG3	2.33	0.63
1:C:105:PRO:HD3	1:C:203:ASN:HB3	1.81	0.62
1:C:47:ALA:O	1:C:218:THR:HG22	2.00	0.62
1:B:46:LEU:HD13	2:B:601:FAD:HM81	1.80	0.62
1:A:372:LEU:HD21	1:A:395:LEU:HD21	1.81	0.61
1:D:47:ALA:O	1:D:218:THR:HG22	2.01	0.61
1:B:92:ILE:HD12	1:B:93:SER:H	1.65	0.61
1:D:142:PHE:CE1	1:D:179:LEU:HD11	2.37	0.60
1:C:69:TYR:CD1	1:C:463:MET:HG3	2.37	0.59
1:C:142:PHE:CE1	1:C:179:LEU:HD11	2.38	0.59
1:A:89:HIS:ND1	1:A:313:LYS:HG2	2.17	0.58
1:A:27:ASN:HB2	1:A:34:TRP:CZ2	2.39	0.58
1:B:142:PHE:CE1	1:B:179:LEU:HD11	2.38	0.58
1:B:27:ASN:HB2	1:B:34:TRP:CZ2	2.39	0.58
1:D:27:ASN:HB2	1:D:34:TRP:HZ2	1.69	0.58
1:A:69:TYR:CG	1:A:463:MET:HG3	2.39	0.58
1:B:92:ILE:O	1:B:93:SER:HB3	2.03	0.57
1:A:62:GLY:HA2	1:A:458:GLN:HE22	1.70	0.56
1:B:69:TYR:CG	1:B:463:MET:HG3	2.39	0.56
1:D:147:VAL:HG21	1:D:160:ARG:NH2	2.20	0.56
1:A:327:ARG:NE	1:A:373:GLU:OE2	2.31	0.56
1:C:27:ASN:HB2	1:C:34:TRP:HZ2	1.70	0.56
1:C:447:ARG:HB2	2:C:601:FAD:H5'1	1.87	0.55
1:C:286:LEU:O	1:C:289:GLN:HB2	2.07	0.55
1:C:104:TYR:CG	1:C:105:PRO:HA	2.42	0.55
1:A:21:GLY:HA2	1:A:462:PHE:CE2	2.43	0.54
1:A:325:PHE:CD1	1:A:372:LEU:CD2	2.91	0.54
1:B:325:PHE:CE1	1:B:372:LEU:HD12	2.42	0.54
1:A:412:HIS:C	1:A:412:HIS:CD2	2.81	0.54
1:D:286:LEU:O	1:D:289:GLN:HB2	2.07	0.54
1:B:334:TYR:N	1:B:334:TYR:CD1	2.76	0.54
1:C:447:ARG:NH1	4:C:603:SO4:O2	2.41	0.53
1:C:22:ALA:O	1:C:469:VAL:HG21	2.08	0.53
1:C:239:LYS:O	1:C:254:LEU:HD13	2.08	0.53
1:D:66:PHE:HB2	1:D:91:ARG:NH2	2.23	0.53
1:A:334:TYR:CD1	1:A:334:TYR:N	2.77	0.52
1:B:11:LEU:O	1:B:265:LEU:HD12	2.09	0.52
1:D:239:LYS:O	1:D:254:LEU:HD13	2.09	0.52
1:A:130:LEU:HD23	1:D:130:LEU:HD23	1.92	0.52
1:B:149:MET:SD	1:B:184:ALA:HA	2.50	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:GLN:HB2	1:A:340:PRO:CD	2.40	0.52
1:D:420:PRO:HD3	1:D:448:PHE:CZ	2.45	0.52
1:D:27:ASN:HB2	1:D:34:TRP:CZ2	2.46	0.51
1:A:334:TYR:HD1	1:A:334:TYR:N	2.08	0.51
1:C:27:ASN:HB2	1:C:34:TRP:CZ2	2.45	0.51
1:A:142:PHE:CE1	1:A:179:LEU:HD11	2.46	0.51
1:D:17:PRO:HD2	2:D:601:FAD:O5'	2.11	0.51
1:B:339:GLN:HB2	1:B:340:PRO:CD	2.40	0.51
1:C:212:PHE:CG	1:C:213:PRO:HD2	2.46	0.50
1:B:334:TYR:N	1:B:334:TYR:HD1	2.08	0.50
1:D:69:TYR:CG	1:D:463:MET:HG3	2.46	0.50
1:D:22:ALA:O	1:D:469:VAL:HG21	2.12	0.50
1:B:21:GLY:HA2	1:B:462:PHE:CE2	2.46	0.50
1:B:412:HIS:C	1:B:412:HIS:CD2	2.82	0.50
1:D:313:LYS:HD3	1:D:316:LEU:HD21	1.94	0.50
1:A:457:ASN:HB2	2:A:601:FAD:O2	2.11	0.49
1:B:89:HIS:ND1	1:B:313:LYS:HG2	2.27	0.49
1:C:79:ALA:HB1	1:C:224:ALA:CB	2.42	0.49
1:B:483:TYR:N	1:B:484:PRO:HD3	2.27	0.49
1:C:69:TYR:CG	1:C:463:MET:HG3	2.47	0.49
1:B:313:LYS:HD2	1:B:313:LYS:N	2.26	0.49
1:A:11:LEU:O	1:A:265:LEU:HD12	2.13	0.49
1:B:306:ARG:HG3	1:B:307:PRO:HD2	1.94	0.49
1:B:447:ARG:HB2	2:B:601:FAD:H5'2	1.95	0.49
1:D:104:TYR:CG	1:D:105:PRO:HA	2.47	0.49
1:B:457:ASN:HB2	2:B:601:FAD:O2	2.12	0.48
1:C:162:TYR:O	1:C:166:VAL:HG23	2.12	0.48
1:D:133:ARG:HG3	1:D:133:ARG:O	2.13	0.48
1:C:147:VAL:HG21	1:C:160:ARG:NH2	2.29	0.48
1:C:57:LEU:HD12	1:C:57:LEU:N	2.28	0.48
1:A:32:PRO:HB3	1:C:504:PHE:CD2	2.49	0.48
1:B:376:GLU:HB2	1:B:382:VAL:CG2	2.43	0.48
1:D:21:GLY:HA2	1:D:462:PHE:CE2	2.49	0.48
1:D:79:ALA:HB1	1:D:224:ALA:CB	2.44	0.47
1:A:149:MET:SD	1:A:184:ALA:HA	2.54	0.47
1:D:483:TYR:N	1:D:484:PRO:HD3	2.30	0.47
1:A:145:TRP:CD2	1:A:179:LEU:HD13	2.49	0.47
1:A:483:TYR:N	1:A:484:PRO:HD3	2.28	0.47
2:B:601:FAD:O4'	2:B:601:FAD:O2'	2.30	0.47
1:D:212:PHE:CG	1:D:213:PRO:HD2	2.50	0.47
2:D:601:FAD:H1'1	2:D:601:FAD:H9	1.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:60:VAL:HG12	1:D:60:VAL:O	2.15	0.46
1:B:188:LEU:O	1:B:192:THR:HG23	2.16	0.46
1:D:57:LEU:HD12	1:D:57:LEU:N	2.30	0.46
1:A:107:GLN:OE1	1:A:107:GLN:N	2.42	0.46
1:A:46:LEU:HD13	2:A:601:FAD:HM81	1.96	0.46
2:A:601:FAD:H1'1	2:A:601:FAD:H9	1.43	0.46
1:B:142:PHE:CD1	1:B:179:LEU:HD11	2.51	0.46
1:B:372:LEU:HD21	1:B:395:LEU:HD11	1.97	0.46
1:C:420:PRO:HD3	1:C:448:PHE:CZ	2.51	0.46
1:C:470:ASP:HB3	1:C:476:ALA:HB3	1.98	0.46
1:B:307:PRO:HG3	1:B:368:TRP:CH2	2.50	0.46
1:B:359:SER:OG	1:B:361:GLU:HB3	2.16	0.46
1:A:431:THR:O	1:A:435:PRO:HG2	2.17	0.45
2:C:601:FAD:N1	2:C:601:FAD:H2'	2.31	0.45
1:B:327:ARG:NH1	4:B:605:SO4:S	2.90	0.45
1:A:307:PRO:HG2	1:A:368:TRP:CZ2	2.52	0.45
1:C:325:PHE:HA	1:C:374:VAL:HG22	1.99	0.45
1:B:92:ILE:HG21	1:B:312:ASP:OD1	2.16	0.45
1:D:276:GLU:OE1	1:D:276:GLU:HA	2.16	0.45
1:A:376:GLU:HB2	1:A:382:VAL:CG2	2.47	0.45
1:B:412:HIS:HD2	1:B:413:ARG:N	2.13	0.45
1:B:104:TYR:CD2	1:B:104:TYR:C	2.90	0.45
1:D:414:ARG:HH11	1:D:414:ARG:HG2	1.82	0.45
1:B:162:TYR:O	1:B:166:VAL:HG23	2.17	0.45
1:A:412:HIS:HD2	1:A:413:ARG:N	2.15	0.45
1:A:62:GLY:HA2	1:A:458:GLN:NE2	2.31	0.45
1:B:494:GLU:HG2	1:D:477:VAL:HA	1.98	0.45
2:C:601:FAD:H1'1	2:C:601:FAD:H9	1.52	0.45
1:D:162:TYR:O	1:D:166:VAL:HG23	2.17	0.45
1:A:242:VAL:HG13	1:A:252:VAL:HG13	1.98	0.44
1:A:59:ASP:HB3	1:A:62:GLY:H	1.82	0.44
1:B:313:LYS:HG3	1:B:333:ASN:HB3	1.99	0.44
1:B:130:LEU:HD23	1:C:130:LEU:HD23	1.97	0.44
1:C:21:GLY:HA2	1:C:462:PHE:CE2	2.52	0.44
1:B:145:TRP:CD2	1:B:179:LEU:HD13	2.53	0.44
1:B:63:HIS:NE2	1:B:334:TYR:HE2	2.15	0.44
1:D:60:VAL:HG13	1:D:415:PHE:CE2	2.52	0.44
1:D:430:LEU:HA	1:D:430:LEU:HD12	1.74	0.44
1:B:372:LEU:HD11	1:B:395:LEU:HD21	2.00	0.44
1:C:414:ARG:HH11	1:C:414:ARG:HG2	1.82	0.44
1:B:61:GLY:O	2:B:601:FAD:C10	2.66	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:133:ARG:O	1:C:133:ARG:HG3	2.17	0.44
1:B:290:LEU:HD13	1:B:451:TRP:CD2	2.53	0.44
1:C:142:PHE:HA	1:C:176:CYS:HB3	1.99	0.43
1:C:483:TYR:N	1:C:484:PRO:HD3	2.33	0.43
1:A:89:HIS:CE1	1:A:313:LYS:HG2	2.52	0.43
1:C:18:THR:OG1	1:C:461:SER:HB3	2.18	0.43
1:D:412:HIS:CD2	1:D:412:HIS:C	2.90	0.43
1:D:69:TYR:CE1	1:D:463:MET:HG3	2.53	0.43
1:B:142:PHE:HA	1:B:176:CYS:HB3	1.99	0.43
1:C:79:ALA:HB1	1:C:224:ALA:HB1	2.00	0.43
1:A:142:PHE:HA	1:A:176:CYS:HB3	2.00	0.43
1:B:245:VAL:O	1:B:278:MET:HA	2.17	0.43
1:D:332:SER:HB3	1:D:369:SER:N	2.33	0.43
1:B:69:TYR:CE1	1:B:463:MET:HG3	2.52	0.43
1:C:412:HIS:CD2	1:C:412:HIS:C	2.89	0.43
1:C:447:ARG:HB2	2:C:601:FAD:C5'	2.47	0.43
1:A:306:ARG:HA	1:A:307:PRO:HD3	1.76	0.43
1:D:470:ASP:HB3	1:D:476:ALA:HB3	1.99	0.43
1:A:323:CYS:HB2	1:A:325:PHE:CE1	2.54	0.43
1:A:69:TYR:CE1	1:A:463:MET:HG3	2.54	0.42
1:A:94:TYR:N	1:A:94:TYR:CD1	2.87	0.42
1:C:332:SER:HB3	1:C:369:SER:N	2.34	0.42
1:B:27:ASN:HB2	1:B:34:TRP:HZ2	1.83	0.42
1:A:27:ASN:HB2	1:A:34:TRP:HZ2	1.81	0.42
1:A:10:VAL:HA	1:A:264:LYS:O	2.19	0.42
1:B:432:GLN:O	1:B:436:LYS:HG3	2.18	0.42
1:B:327:ARG:NH1	4:B:605:SO4:O1	2.52	0.42
1:D:105:PRO:HD3	1:D:203:ASN:HB3	2.02	0.42
1:A:245:VAL:O	1:A:278:MET:HA	2.19	0.42
1:A:325:PHE:CE1	1:A:372:LEU:CD2	3.03	0.42
1:D:42:THR:HA	1:D:43:PRO:HD3	1.87	0.42
1:B:51:VAL:HG22	1:B:52:THR:O	2.19	0.42
1:C:430:LEU:HA	1:C:430:LEU:HD12	1.74	0.42
1:D:142:PHE:HA	1:D:176:CYS:HB3	2.02	0.42
1:A:432:GLN:O	1:A:436:LYS:HG3	2.19	0.42
1:B:10:VAL:HA	1:B:264:LYS:O	2.20	0.42
1:B:167:TRP:HH2	1:B:453:TYR:CD1	2.38	0.42
1:D:18:THR:OG1	1:D:461:SER:HB3	2.20	0.42
1:D:24:LYS:HA	1:D:24:LYS:HD2	1.91	0.42
1:D:325:PHE:CD1	1:D:372:LEU:HD13	2.55	0.42
1:A:372:LEU:HD11	1:A:395:LEU:HD11	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:325:PHE:HA	1:B:374:VAL:HG22	2.01	0.42
1:D:213:PRO:HG3	1:D:221:ILE:HG13	2.01	0.42
1:B:328:ALA:HA	1:B:371:MET:O	2.20	0.41
1:B:504:PHE:CD2	1:D:32:PRO:HB3	2.55	0.41
1:A:453:TYR:CZ	3:A:602:NAI:H42N	2.54	0.41
1:D:142:PHE:CD1	1:D:179:LEU:HD11	2.55	0.41
1:D:3:HIS:HA	1:D:4:PRO:HD3	1.79	0.41
1:A:325:PHE:HA	1:A:374:VAL:HG22	2.03	0.41
1:A:332:SER:HB3	1:A:369:SER:N	2.36	0.41
1:C:142:PHE:CD1	1:C:179:LEU:HD11	2.55	0.41
2:B:601:FAD:H9	2:B:601:FAD:H1'1	1.49	0.41
1:C:212:PHE:CD2	1:C:213:PRO:HD2	2.56	0.41
1:C:332:SER:HB3	1:C:369:SER:H	1.86	0.41
1:A:162:TYR:O	1:A:166:VAL:HG23	2.21	0.41
1:C:413:ARG:HD3	1:C:415:PHE:CZ	2.56	0.41
1:D:141:THR:HG1	1:D:144:GLU:HG3	1.86	0.40
1:D:332:SER:HB3	1:D:369:SER:H	1.86	0.40
1:D:413:ARG:HD3	1:D:415:PHE:CZ	2.56	0.40
1:A:142:PHE:CD1	1:A:179:LEU:HD11	2.56	0.40
1:A:8:VAL:O	1:A:262:TYR:HA	2.22	0.40
1:B:92:ILE:HG23	1:B:93:SER:N	2.36	0.40
1:C:69:TYR:CE1	1:C:463:MET:HG3	2.56	0.40
1:C:66:PHE:HB2	1:C:91:ARG:HH12	1.86	0.40
1:B:472:ILE:O	1:D:500:GLY:HA3	2.22	0.40
1:B:483:TYR:N	1:B:484:PRO:CD	2.84	0.40
1:D:20:LEU:HD23	1:D:20:LEU:HA	1.92	0.40
1:B:313:LYS:CD	1:B:313:LYS:N	2.85	0.40
1:C:445:ARG:NE	1:C:481:LEU:HD22	2.36	0.40
1:A:494:GLU:HG2	1:C:477:VAL:HA	2.03	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	488/513 (95%)	469 (96%)	19 (4%)	0	100	100
1	B	488/513 (95%)	468 (96%)	20 (4%)	0	100	100
1	C	495/513 (96%)	478 (97%)	17 (3%)	0	100	100
1	D	498/513 (97%)	482 (97%)	16 (3%)	0	100	100
All	All	1969/2052 (96%)	1897 (96%)	72 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	386/432 (89%)	376 (97%)	10 (3%)	54	84
1	B	389/432 (90%)	378 (97%)	11 (3%)	51	82
1	C	383/432 (89%)	374 (98%)	9 (2%)	58	87
1	D	384/432 (89%)	374 (97%)	10 (3%)	54	84
All	All	1542/1728 (89%)	1502 (97%)	40 (3%)	54	84

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

Mol	Chain	Res	Type
1	A	104	TYR
1	A	149	MET
1	A	169	VAL
1	A	175	GLN
1	A	334	TYR
1	A	384	GLN
1	A	417	HIS
1	A	430	LEU
1	A	470	ASP
1	A	485	ASP
1	B	92	ILE
1	B	104	TYR
1	B	149	MET

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Mol	Chain	Res	Type
1	B	169	VAL
1	B	175	GLN
1	B	334	TYR
1	B	372	LEU
1	B	417	HIS
1	B	430	LEU
1	B	470	ASP
1	B	485	ASP
1	C	7	SER
1	C	27	ASN
1	C	175	GLN
1	C	218	THR
1	C	360	THR
1	C	430	LEU
1	C	470	ASP
1	C	485	ASP
1	C	487	VAL
1	D	7	SER
1	D	27	ASN
1	D	84	ASP
1	D	175	GLN
1	D	218	THR
1	D	360	THR
1	D	430	LEU
1	D	470	ASP
1	D	485	ASP
1	D	487	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	GLN
1	A	458	GLN
1	B	98	GLN
1	B	458	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

16 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	-	52,58,58	2.01	19 (36%)	52,89,89	2.76	21 (40%)
3	NAI	A	602	-	41,48,48	1.76	15 (36%)	46,73,73	2.36	7 (15%)
4	SO4	A	603	-	4,4,4	0.11	0	6,6,6	0.12	0
4	SO4	A	604	-	4,4,4	0.23	0	6,6,6	0.24	0
2	FAD	B	601	-	52,58,58	2.06	19 (36%)	52,89,89	2.62	20 (38%)
3	NAI	B	602	-	41,48,48	1.82	14 (34%)	46,73,73	2.27	9 (19%)
4	SO4	B	603	-	4,4,4	0.17	0	6,6,6	0.29	0
4	SO4	B	604	-	4,4,4	0.18	0	6,6,6	0.27	0
4	SO4	B	605	-	4,4,4	0.29	0	6,6,6	0.17	0
2	FAD	C	601	-	52,58,58	1.92	17 (32%)	52,89,89	2.48	15 (28%)
4	SO4	C	602	-	4,4,4	0.14	0	6,6,6	0.23	0
4	SO4	C	603	-	4,4,4	0.26	0	6,6,6	0.18	0
2	FAD	D	601	-	52,58,58	2.03	19 (36%)	52,89,89	2.60	18 (34%)
4	SO4	D	602	-	4,4,4	0.32	0	6,6,6	0.31	0
4	SO4	D	603	-	4,4,4	0.17	0	6,6,6	0.23	0
4	SO4	D	604	-	4,4,4	0.25	0	6,6,6	0.08	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FAD	A	601	-	-	0/30/50/50	0/6/6/6
3	NAI	A	602	-	-	0/25/72/72	0/5/5/5
4	SO4	A	603	-	-	0/0/0/0	0/0/0/0
4	SO4	A	604	-	-	0/0/0/0	0/0/0/0
2	FAD	B	601	-	-	0/30/50/50	0/6/6/6
3	NAI	B	602	-	-	0/25/72/72	0/5/5/5
4	SO4	B	603	-	-	0/0/0/0	0/0/0/0
4	SO4	B	604	-	-	0/0/0/0	0/0/0/0
4	SO4	B	605	-	-	0/0/0/0	0/0/0/0
2	FAD	C	601	-	-	0/30/50/50	0/6/6/6
4	SO4	C	602	-	-	0/0/0/0	0/0/0/0
4	SO4	C	603	-	-	0/0/0/0	0/0/0/0
2	FAD	D	601	-	-	0/30/50/50	0/6/6/6
4	SO4	D	602	-	-	0/0/0/0	0/0/0/0
4	SO4	D	603	-	-	0/0/0/0	0/0/0/0
4	SO4	D	604	-	-	0/0/0/0	0/0/0/0

All (103) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	FAD	C2B-C1B	-3.80	1.47	1.53
2	A	601	FAD	O2'-C2'	-3.42	1.35	1.43
3	B	602	NAI	C4N-C5N	-3.38	1.41	1.49
2	A	601	FAD	C2B-C1B	-3.16	1.48	1.53
3	A	602	NAI	C4N-C5N	-3.15	1.42	1.49
2	D	601	FAD	C1'-N10	-3.15	1.45	1.48
2	C	601	FAD	C2B-C1B	-3.08	1.48	1.53
2	C	601	FAD	C1'-N10	-3.04	1.45	1.48
2	B	601	FAD	O2'-C2'	-3.01	1.36	1.43
2	D	601	FAD	C2B-C1B	-3.01	1.48	1.53
2	D	601	FAD	PA-O5B	-2.94	1.46	1.59
2	A	601	FAD	C1'-N10	-2.84	1.45	1.48
2	A	601	FAD	O4'-C4'	-2.74	1.37	1.43
2	D	601	FAD	O2'-C2'	-2.72	1.37	1.43
2	B	601	FAD	PA-O5B	-2.71	1.47	1.59
2	B	601	FAD	O4'-C4'	-2.68	1.37	1.43
3	A	602	NAI	C2B-C1B	-2.63	1.49	1.53
2	B	601	FAD	C1'-N10	-2.61	1.45	1.48
2	C	601	FAD	O4'-C4'	-2.59	1.37	1.43
2	A	601	FAD	PA-O5B	-2.58	1.48	1.59
3	B	602	NAI	PA-O2A	-2.57	1.44	1.55
3	B	602	NAI	O2D-C2D	-2.53	1.37	1.43
2	C	601	FAD	O2'-C2'	-2.52	1.37	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	601	FAD	O4'-C4'	-2.51	1.37	1.43
2	C	601	FAD	PA-O5B	-2.45	1.48	1.59
3	B	602	NAI	C2B-C1B	-2.44	1.49	1.53
3	A	602	NAI	O2D-C2D	-2.41	1.37	1.43
2	B	601	FAD	PA-O2A	-2.41	1.44	1.55
2	B	601	FAD	C2B-C3B	-2.37	1.47	1.53
3	B	602	NAI	C2B-C3B	-2.36	1.47	1.53
3	B	602	NAI	PA-O5B	-2.31	1.49	1.59
2	A	601	FAD	O2B-C2B	-2.28	1.37	1.43
3	A	602	NAI	C2B-C3B	-2.26	1.47	1.53
2	B	601	FAD	O3B-C3B	-2.25	1.37	1.43
3	B	602	NAI	O2B-C2B	-2.25	1.37	1.43
2	A	601	FAD	PA-O2A	-2.23	1.45	1.55
2	A	601	FAD	C2B-C3B	-2.23	1.47	1.53
3	A	602	NAI	O2B-C2B	-2.22	1.37	1.43
2	D	601	FAD	C2B-C3B	-2.19	1.47	1.53
3	A	602	NAI	PA-O5B	-2.13	1.50	1.59
2	D	601	FAD	O4B-C4B	-2.13	1.40	1.45
2	A	601	FAD	C4'-C3'	-2.13	1.49	1.53
2	C	601	FAD	O3'-C3'	-2.11	1.38	1.43
2	D	601	FAD	PA-O2A	-2.09	1.46	1.55
2	B	601	FAD	O2B-C2B	-2.08	1.38	1.43
3	A	602	NAI	PA-O2A	-2.07	1.46	1.55
2	A	601	FAD	O3'-C3'	-2.05	1.38	1.43
3	A	602	NAI	O3B-C3B	-2.05	1.38	1.43
2	C	601	FAD	O2B-C2B	-2.02	1.38	1.43
2	D	601	FAD	C10-N10	2.01	1.41	1.39
2	C	601	FAD	C4X-C10	2.05	1.44	1.40
3	B	602	NAI	C2N-C3N	2.12	1.40	1.34
2	A	601	FAD	C5X-N5	2.13	1.38	1.35
2	A	601	FAD	C4X-C10	2.14	1.44	1.40
2	C	601	FAD	C5X-N5	2.15	1.38	1.35
3	A	602	NAI	C6N-N1N	2.17	1.42	1.37
3	B	602	NAI	C6N-N1N	2.22	1.43	1.37
2	B	601	FAD	P-O1P	2.24	1.59	1.51
2	D	601	FAD	P-O1P	2.27	1.59	1.51
2	B	601	FAD	C2A-N3A	2.28	1.36	1.32
2	B	601	FAD	C4X-C10	2.30	1.45	1.40
3	A	602	NAI	C2N-C3N	2.33	1.41	1.34
2	D	601	FAD	C4X-C10	2.33	1.45	1.40
3	B	602	NAI	C2A-N3A	2.35	1.36	1.32
2	A	601	FAD	C2A-N3A	2.38	1.36	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	FAD	C6A-N6A	2.38	1.43	1.34
3	A	602	NAI	C6A-N6A	2.39	1.43	1.34
2	A	601	FAD	C6A-N6A	2.40	1.43	1.34
3	A	602	NAI	C2A-N3A	2.43	1.36	1.32
3	B	602	NAI	C2N-N1N	2.44	1.41	1.37
2	C	601	FAD	P-O1P	2.53	1.60	1.51
2	D	601	FAD	C6A-N6A	2.54	1.44	1.34
2	C	601	FAD	C6A-N6A	2.55	1.44	1.34
3	A	602	NAI	C2N-N1N	2.60	1.42	1.37
2	D	601	FAD	C2-N1	2.65	1.43	1.38
2	C	601	FAD	C2A-N3A	2.69	1.36	1.32
3	B	602	NAI	C6A-N6A	2.72	1.45	1.34
2	D	601	FAD	C2A-N3A	2.73	1.37	1.32
2	D	601	FAD	C5X-N5	2.76	1.39	1.35
2	B	601	FAD	C2-N1	2.82	1.44	1.38
2	B	601	FAD	C5X-N5	2.84	1.39	1.35
2	A	601	FAD	C2-N1	2.96	1.44	1.38
2	C	601	FAD	C2-N1	3.09	1.44	1.38
2	C	601	FAD	C4X-N5	3.18	1.38	1.33
2	C	601	FAD	C4-N3	3.38	1.39	1.33
2	D	601	FAD	C4-N3	3.44	1.39	1.33
2	A	601	FAD	C4-N3	3.51	1.39	1.33
3	A	602	NAI	C7N-N7N	3.54	1.43	1.33
2	D	601	FAD	C4X-N5	3.69	1.39	1.33
3	B	602	NAI	C7N-N7N	3.73	1.44	1.33
2	C	601	FAD	C4-C4X	3.78	1.49	1.41
2	B	601	FAD	C4-N3	3.83	1.39	1.33
2	A	601	FAD	C4X-N5	3.92	1.39	1.33
3	A	602	NAI	C6N-C5N	4.00	1.40	1.33
2	B	601	FAD	C4X-N5	4.01	1.39	1.33
3	B	602	NAI	C6N-C5N	4.03	1.40	1.33
2	B	601	FAD	C4-C4X	4.41	1.50	1.41
2	A	601	FAD	C4-C4X	4.55	1.50	1.41
2	D	601	FAD	C4-C4X	4.79	1.51	1.41
2	C	601	FAD	O4-C4	5.26	1.37	1.24
2	D	601	FAD	O4-C4	5.53	1.38	1.24
2	A	601	FAD	O4-C4	5.68	1.39	1.24
2	B	601	FAD	O4-C4	5.70	1.39	1.24

All (90) bond angle outliers are listed below:

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	NAI	N3A-C2A-N1A	-12.02	119.43	128.87
3	B	602	NAI	N3A-C2A-N1A	-12.00	119.45	128.87
2	A	601	FAD	N3A-C2A-N1A	-10.75	120.43	128.87
2	B	601	FAD	N3A-C2A-N1A	-10.42	120.69	128.87
2	D	601	FAD	N3A-C2A-N1A	-10.22	120.84	128.87
2	C	601	FAD	N3A-C2A-N1A	-10.17	120.89	128.87
2	D	601	FAD	O2P-P-O3P	-6.07	79.28	105.27
2	A	601	FAD	C1'-N10-C9A	-5.72	112.19	118.83
2	A	601	FAD	C4B-O4B-C1B	-4.74	104.62	109.64
2	B	601	FAD	N3-C2-N1	-4.20	120.62	127.69
2	C	601	FAD	O2P-P-O5'	-4.16	88.39	108.24
2	C	601	FAD	N3-C2-N1	-4.06	120.85	127.69
3	A	602	NAI	C1B-N9A-C4A	-4.05	122.28	126.81
2	A	601	FAD	N3-C2-N1	-3.91	121.11	127.69
2	B	601	FAD	C1'-N10-C9A	-3.90	114.30	118.83
2	D	601	FAD	N3-C2-N1	-3.88	121.16	127.69
2	C	601	FAD	C4X-C10-N10	-3.70	117.83	120.52
2	D	601	FAD	C1'-N10-C9A	-3.51	114.76	118.83
2	A	601	FAD	C4X-C10-N10	-3.47	118.00	120.52
2	B	601	FAD	C4X-C4-N3	-3.40	119.08	123.52
2	A	601	FAD	C4X-C4-N3	-3.39	119.09	123.52
2	D	601	FAD	C4X-C4-N3	-3.29	119.22	123.52
2	B	601	FAD	O2P-P-O3P	-3.29	91.16	105.27
3	A	602	NAI	C4B-O4B-C1B	-3.26	106.19	109.64
2	D	601	FAD	C4-C4X-C10	-3.14	117.93	119.94
2	B	601	FAD	C4-C4X-C10	-3.10	117.96	119.94
2	C	601	FAD	C4X-C4-N3	-2.98	119.62	123.52
2	A	601	FAD	O2P-P-O3P	-2.94	92.67	105.27
2	C	601	FAD	C4-C4X-C10	-2.91	118.08	119.94
3	A	602	NAI	O2A-PA-O3	-2.85	93.03	105.27
3	B	602	NAI	C4B-O4B-C1B	-2.82	106.65	109.64
3	B	602	NAI	O2A-PA-O5B	-2.81	94.83	108.24
2	A	601	FAD	O2P-P-O5'	-2.79	94.93	108.24
2	A	601	FAD	O2A-PA-O3P	-2.69	93.74	105.27
2	B	601	FAD	C4X-C10-N10	-2.68	118.58	120.52
2	C	601	FAD	C1'-N10-C9A	-2.66	115.74	118.83
3	B	602	NAI	C1B-N9A-C4A	-2.65	123.85	126.81
2	A	601	FAD	O2A-PA-O5B	-2.59	95.91	108.24
2	A	601	FAD	C4-C4X-C10	-2.55	118.31	119.94
2	B	601	FAD	O2A-PA-O3P	-2.44	94.81	105.27
3	B	602	NAI	C3N-C2N-N1N	-2.43	119.67	123.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	FAD	O2'-C2'-C1'	-2.41	103.99	109.93
2	B	601	FAD	O2P-P-O5'	-2.39	96.82	108.24
2	A	601	FAD	C1'-C2'-C3'	-2.33	103.17	109.82
2	C	601	FAD	O2P-P-O3P	-2.23	95.71	105.27
2	D	601	FAD	C5B-C4B-C3B	-2.13	106.95	115.20
2	D	601	FAD	O2A-PA-O5B	-2.13	98.10	108.24
2	D	601	FAD	C4B-O4B-C1B	-2.11	107.41	109.64
2	B	601	FAD	C4B-O4B-C1B	-2.06	107.46	109.64
2	D	601	FAD	O3'-C3'-C2'	2.03	114.00	108.73
2	B	601	FAD	O5B-PA-O1A	2.06	117.66	109.21
2	A	601	FAD	C6-C5X-N5	2.12	121.56	118.92
2	A	601	FAD	O5B-C5B-C4B	2.19	117.01	109.09
2	D	601	FAD	O2P-P-O1P	2.24	124.21	112.56
3	B	602	NAI	O5B-PA-O1A	2.25	118.42	109.21
2	C	601	FAD	O5B-PA-O1A	2.25	118.43	109.21
2	B	601	FAD	O5'-P-O1P	2.33	118.73	109.21
2	D	601	FAD	O5B-PA-O1A	2.35	118.85	109.21
2	B	601	FAD	C6-C5X-N5	2.42	121.93	118.92
2	A	601	FAD	O5'-P-O1P	2.42	119.12	109.21
2	D	601	FAD	C5X-C9A-N10	2.43	119.40	117.58
2	C	601	FAD	C4-C4X-N5	2.50	121.74	118.70
2	D	601	FAD	O2A-PA-O1A	2.52	125.66	112.56
3	B	602	NAI	O4B-C1B-N9A	2.65	113.11	108.11
3	B	602	NAI	C5N-C4N-C3N	2.70	119.40	112.41
3	A	602	NAI	C5N-C4N-C3N	2.73	119.48	112.41
2	D	601	FAD	O4B-C1B-N9A	2.82	113.43	108.11
2	A	601	FAD	O5B-PA-O1A	2.88	121.01	109.21
2	B	601	FAD	O3'-C3'-C2'	2.98	116.46	108.73
2	C	601	FAD	O5'-P-O1P	3.00	121.50	109.21
2	A	601	FAD	C4-C4X-N5	3.09	122.46	118.70
2	C	601	FAD	O4B-C1B-N9A	3.13	114.02	108.11
2	A	601	FAD	C5X-C9A-N10	3.33	120.07	117.58
2	B	601	FAD	C4-C4X-N5	3.34	122.76	118.70
2	B	601	FAD	C5X-C9A-N10	3.37	120.11	117.58
2	C	601	FAD	C5X-C9A-N10	3.51	120.20	117.58
2	B	601	FAD	C4X-N5-C5X	3.52	120.86	116.72
2	C	601	FAD	C4X-N5-C5X	3.61	120.97	116.72
3	A	602	NAI	O4D-C1D-N1N	3.67	115.08	108.09
2	A	601	FAD	O4B-C1B-N9A	3.71	115.12	108.11
2	B	601	FAD	O4B-C1B-N9A	3.77	115.23	108.11
2	D	601	FAD	C4X-N5-C5X	3.96	121.39	116.72
2	A	601	FAD	C4X-N5-C5X	3.99	121.43	116.72

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	D	601	FAD	C4-C4X-N5	4.20	123.81	118.70
3	B	602	NAI	O4D-C1D-N1N	4.23	116.17	108.09
3	A	602	NAI	O4B-C1B-N9A	4.33	116.29	108.11
2	D	601	FAD	C4-N3-C2	6.56	120.63	115.16
2	A	601	FAD	C4-N3-C2	6.67	120.73	115.16
2	C	601	FAD	C4-N3-C2	6.79	120.82	115.16
2	B	601	FAD	C4-N3-C2	6.96	120.96	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 24 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	FAD	3	0
3	A	602	NAI	2	0
2	B	601	FAD	6	0
3	B	602	NAI	2	0
4	B	605	SO4	2	0
2	C	601	FAD	5	0
4	C	603	SO4	1	0
2	D	601	FAD	3	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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	494/513 (96%)	0.29	2 (0%) 93 92	25, 46, 76, 103	0
1	B	494/513 (96%)	0.26	3 (0%) 90 88	24, 47, 77, 121	0
1	C	499/513 (97%)	0.37	9 (1%) 71 66	24, 53, 84, 106	0
1	D	502/513 (97%)	0.38	11 (2%) 65 59	23, 54, 86, 104	0
All	All	1989/2052 (96%)	0.32	25 (1%) 79 75	23, 50, 81, 121	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	236	PHE	3.3
1	C	245	VAL	2.9
1	C	273	PHE	2.7
1	B	184	ALA	2.6
1	C	242	VAL	2.6
1	A	62	GLY	2.5
1	B	23	ALA	2.4
1	D	218	THR	2.4
1	B	497	LEU	2.3
1	C	46	LEU	2.3
1	D	240	GLY	2.3
1	C	11	LEU	2.3
1	D	223	ILE	2.2
1	C	411	TYR	2.2
1	D	221	ILE	2.2
1	D	498	VAL	2.2
1	D	411	TYR	2.1
1	C	20	LEU	2.1
1	D	259	THR	2.1
1	D	273	PHE	2.1
1	C	58	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	36	ILE	2.1
1	A	307	PRO	2.0
1	D	66	PHE	2.0
1	C	300	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	SO4	C	603	5/5	0.96	0.27	2.00	46,49,61,87	5
4	SO4	B	605	5/5	0.95	0.23	0.63	70,74,84,89	0
3	NAI	B	602	44/44	0.97	0.21	0.19	29,48,63,71	0
4	SO4	D	602	5/5	0.99	0.24	0.15	30,31,55,82	5
2	FAD	B	601	53/53	0.98	0.20	-0.04	18,43,62,66	0
2	FAD	D	601	53/53	0.95	0.21	-0.26	26,57,91,107	0
2	FAD	A	601	53/53	0.98	0.20	-0.37	25,44,63,69	0
2	FAD	C	601	53/53	0.97	0.20	-0.60	24,60,86,96	0
3	NAI	A	602	44/44	0.96	0.19	-1.10	22,54,71,78	0
4	SO4	A	603	5/5	0.98	0.17	-	29,37,66,74	5
4	SO4	C	602	5/5	0.94	0.16	-	63,74,90,102	0
4	SO4	D	604	5/5	0.95	0.22	-	51,102,114,116	0
4	SO4	A	604	5/5	0.93	0.23	-	53,68,89,90	0
4	SO4	B	604	5/5	0.97	0.21	-	45,66,73,76	5
4	SO4	B	603	5/5	0.93	0.12	-	68,94,101,101	0
4	SO4	D	603	5/5	0.94	0.21	-	80,89,94,94	5

6.5 Other polymers [i](#)

There are no such residues in this entry.