



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 01:16 AM GMT

PDB ID : 3GDP
Title : Hydroxynitrile lyase from almond, monoclinic crystal form
Authors : Dreveny, I.; Gruber, K.; Kratky, C.
Deposited on : 2009-02-24
Resolution : 1.57 Å(reported)

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

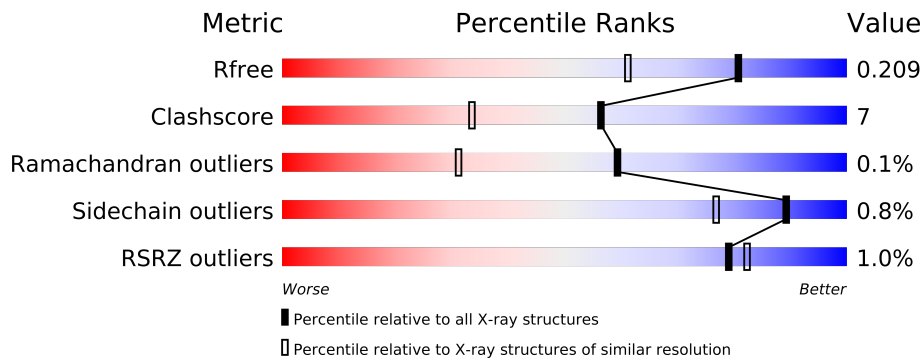
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.57 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}	66092	2778 (1.60-1.56)
Clashscore	79885	3207 (1.60-1.56)
Ramachandran outliers	78287	3107 (1.60-1.56)
Sidechain outliers	78261	3104 (1.60-1.56)
RSRZ outliers	66119	2778 (1.60-1.56)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	521	
1	B	521	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	IPA	A	522	-	X
4	NAG	A	524	-	X
4	NAG	A	533	-	X
9	NDG	B	533	-	X

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 10004 atoms, of which 0 are hydrogen and 0 are deuterium.

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

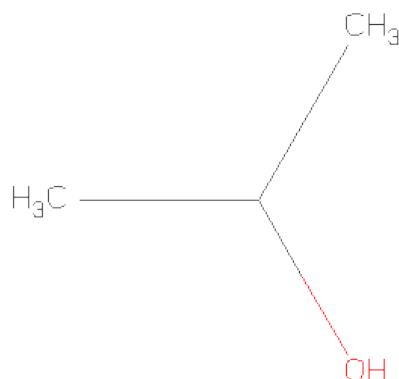
- Molecule 1 is a protein called R-oxynitrile lyase isoenzyme 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	521	Total	C	N	O	S	0	10	0
			4022	2556	669	789	8			
1	B	521	Total	C	N	O	S	0	8	0
			4015	2550	667	790	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	346	SER	ALA	SEE REMARK 999	UNP Q945K2
B	346	SER	ALA	SEE REMARK 999	UNP Q945K2

- Molecule 2 is ISOPROPYL ALCOHOL (three-letter code: IPA) (formula: C₃H₈O).



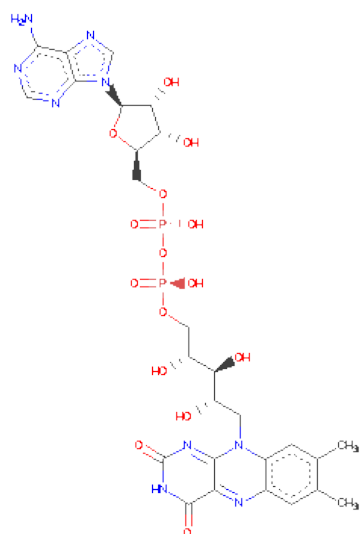
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	3	1		

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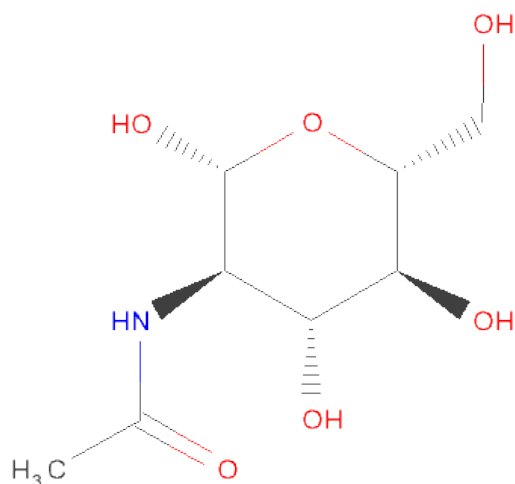
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			4	3	1		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is a polymer of unknown type called SUGAR (5-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	5	Total	C	N	O	0	0
			60	34	2	24		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	346	SER	ALA	SEE REMARK 999	UNP Q945K2

- Molecule 6 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	3	Total	C	N	O	0	0
			38	22	2	14		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	346	SER	ALA	SEE REMARK 999	UNP Q945K2

- Molecule 7 is a polymer of unknown type called SUGAR (5-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	B	5	Total	C	N	O	0	0
			60	34	2	24		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	346	SER	ALA	SEE REMARK 999	UNP Q945K2

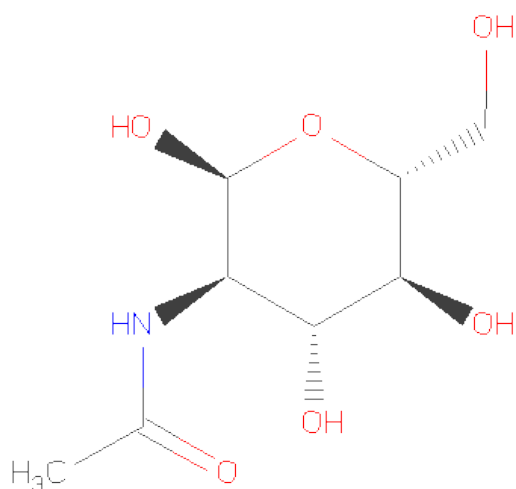
- Molecule 8 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	B	4	Total	C	N	O	0	0
			50	28	2	20		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	346	SER	ALA	SEE REMARK 999	UNP Q945K2

- Molecule 9 is SUGAR (2-(ACETYLAMINO)-2-DEOXY-A-D-GLUCOPYRANOSE) (three-letter code: NDG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 10 is water.

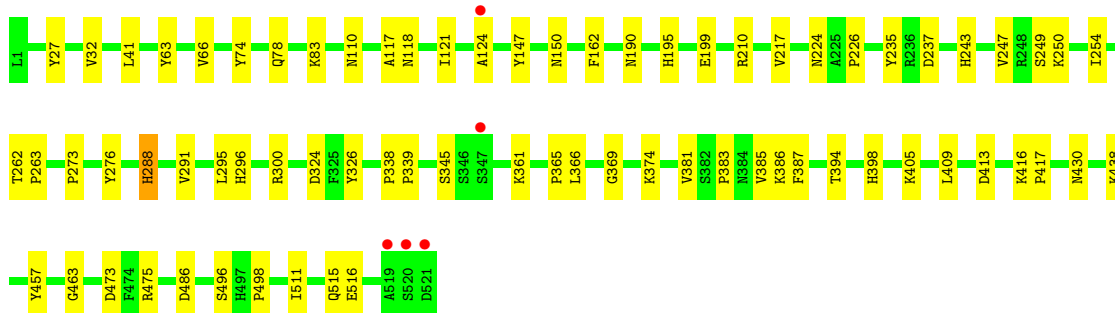
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	823	Total	O	0	0
			823	823		
10	B	766	Total	O	0	0
			766	766		

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.

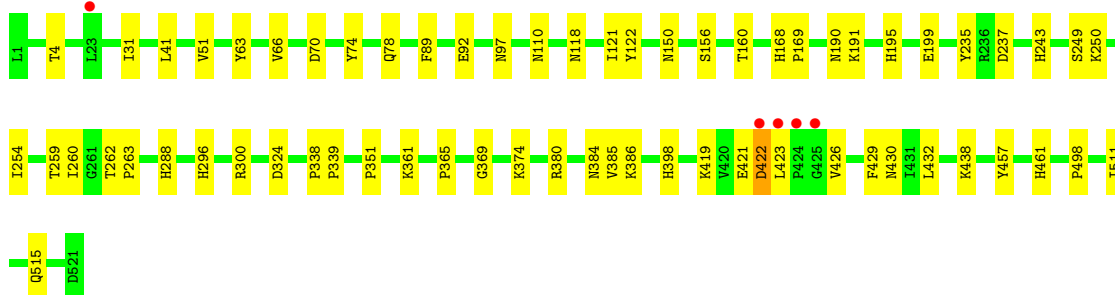
- Molecule 1: R-oxynitrile lyase isoenzyme 1

Chain A: 



- Molecule 1: R-oxynitrile lyase isoenzyme 1

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	69.05Å 93.71Å 87.26Å 90.00° 106.39° 90.00°	Depositor
Resolution (Å)	38.25 – 1.57 38.25 – 1.57	Depositor EDS
% Data completeness (in resolution range)	95.6 (38.25-1.57) 95.7 (38.25-1.57)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 1.57Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.186 , 0.210 0.185 , 0.209	Depositor DCC
R_{free} test set	14166 reflections (11.07%)	DCC
Wilson B-factor (Å ²)	11.7	Xtriage
Anisotropy	0.663	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 142084 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10004	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 48.30 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 8.7740e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NAG, NDG, IPA, FUC, FUL, FAD, MAN

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	0.31	0/4181	0.63	0/5702
1	B	0.32	0/4161	0.63	0/5675
All	All	0.31	0/8342	0.63	0/11377

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4022	0	3882	61	0
1	B	4015	0	3874	45	0
2	A	4	0	8	1	0
2	B	4	0	8	0	0
3	A	53	0	31	2	0
3	B	53	0	31	1	0
4	A	28	0	26	0	0
4	B	14	0	13	0	0
5	A	60	0	52	1	0
6	A	38	0	34	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	60	0	52	1	0
8	B	50	0	43	0	0
9	B	14	0	13	0	0
10	A	823	0	0	9	0
10	B	766	0	0	5	0
All	All	10004	0	8067	108	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (108) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:295[B]:LEU:CD2	1:A:385[B]:VAL:CG2	2.03	1.35
1:A:475[B]:ARG:NH1	10:A:1007:HOH:O	1.60	1.31
1:A:295[B]:LEU:HD21	1:A:385[B]:VAL:HG21	1.19	1.19
1:A:295[B]:LEU:CD2	1:A:385[B]:VAL:HG23	1.75	1.06
1:A:295[B]:LEU:HD22	1:A:385[B]:VAL:HG23	1.42	1.01
1:A:295[B]:LEU:HD23	1:A:385[B]:VAL:CG2	1.92	0.98
1:A:295[B]:LEU:HD21	1:A:385[B]:VAL:CG2	1.78	0.97
1:A:150:ASN:ND2	1:A:190:ASN:HD22	1.65	0.92
1:A:210[B]:ARG:NH1	10:A:959:HOH:O	2.10	0.79
1:B:421:GLU:HG2	1:B:423:LEU:HG	1.67	0.76
1:A:295[B]:LEU:HD23	1:A:385[B]:VAL:HG22	1.65	0.76
1:B:70[B]:ASP:OD1	10:B:1555:HOH:O	2.02	0.76
1:A:121:ILE:HD12	2:A:522:IPA:H2	1.70	0.72
1:A:345:SER:HB3	10:A:647:HOH:O	1.89	0.71
1:B:118:ASN:O	1:B:121:ILE:HG13	1.93	0.69
1:B:374:LYS:HD3	1:B:384:ASN:HD21	1.60	0.67
1:A:295[B]:LEU:HG	1:A:387:PHE:HE2	1.59	0.67
1:A:413:ASP:HA	1:A:416:LYS:HD2	1.77	0.66
1:B:422:ASP:O	1:B:423:LEU:HD23	1.95	0.66
1:A:374:LYS:HD3	1:A:381:VAL:CG1	2.27	0.65
5:A:526:NAG:H61	5:A:527:FUL:O2	1.97	0.64
1:B:237:ASP:OD2	1:B:243:HIS:HE1	1.82	0.62
1:A:511:ILE:O	1:A:515:GLN:HG3	1.99	0.62
1:A:150:ASN:HD21	1:A:190:ASN:HD22	1.46	0.62
1:A:237:ASP:OD2	1:A:243:HIS:HE1	1.83	0.62
1:B:511:ILE:O	1:B:515:GLN:HG3	2.01	0.60
1:A:118:ASN:O	1:A:121:ILE:HG12	2.02	0.60
1:B:63:TYR:O	1:B:66:VAL:HG22	2.02	0.60
1:A:475[A]:ARG:NH1	10:A:544:HOH:O	2.35	0.59
1:A:295[B]:LEU:HG	1:A:387:PHE:CE2	2.38	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:31:ILE:N	1:B:31:ILE:HD12	2.19	0.58
1:B:374:LYS:HD3	1:B:384:ASN:ND2	2.21	0.55
7:B:524:NAG:O6	7:B:526:FUC:H3	2.07	0.54
1:B:369:GLY:HA3	1:B:386:LYS:O	2.08	0.54
1:A:117:ALA:HB1	1:A:121:ILE:HD11	1.90	0.53
1:A:374:LYS:HD3	1:A:381:VAL:HG11	1.91	0.53
1:B:374:LYS:HE2	10:B:1111:HOH:O	2.10	0.52
1:A:83:LYS:HD2	10:A:799:HOH:O	2.10	0.52
1:B:438:LYS:HE2	10:B:1151:HOH:O	2.10	0.51
1:A:369:GLY:HA3	1:A:386:LYS:O	2.11	0.51
1:B:41:LEU:HD21	1:B:254:ILE:HG21	1.92	0.51
1:A:262:THR:HB	1:A:263:PRO:HD3	1.92	0.50
1:A:124:ALA:HB1	10:A:611:HOH:O	2.12	0.50
1:A:63:TYR:O	1:A:66:VAL:HG22	2.11	0.50
1:B:426:VAL:HG22	10:B:981:HOH:O	2.12	0.50
1:B:121:ILE:C	1:B:121:ILE:HD12	2.32	0.50
1:A:416:LYS:HB2	1:A:417:PRO:HD3	1.94	0.49
1:B:121:ILE:HD12	1:B:122:TYR:N	2.27	0.49
1:A:235:TYR:CZ	1:A:243:HIS:HB2	2.48	0.48
1:A:475[B]:ARG:NH1	10:A:1006:HOH:O	2.45	0.48
1:A:475[A]:ARG:NH2	1:A:516:GLU:OE2	2.46	0.48
1:A:405:LYS:O	1:A:409[A]:LEU:HD13	2.13	0.48
1:B:300:ARG:HB2	1:B:457:TYR:CD1	2.50	0.47
1:A:438:LYS:HG3	10:A:991:HOH:O	2.13	0.47
1:B:262:THR:HB	1:B:263:PRO:HD3	1.97	0.47
1:A:110:ASN:HB2	3:A:523:FAD:C5X	2.44	0.47
1:A:296:HIS:HB3	1:A:365:PRO:HD2	1.96	0.47
1:A:41:LEU:HD21	1:A:254:ILE:HG21	1.98	0.47
1:B:31:ILE:HD13	1:B:51:VAL:HG13	1.96	0.46
1:A:300:ARG:HB2	1:A:457:TYR:CD1	2.51	0.46
1:B:235:TYR:CZ	1:B:243:HIS:HB2	2.51	0.46
1:A:224:ASN:HB3	1:A:226:PRO:O	2.16	0.46
1:B:110:ASN:HB2	3:B:523:FAD:C5X	2.45	0.46
1:B:419:LYS:HE2	1:B:429:PHE:CE1	2.51	0.46
1:B:150:ASN:HD21	1:B:190:ASN:HB3	1.81	0.45
1:A:473:ASP:HB2	1:A:475[A]:ARG:NH1	2.31	0.45
1:A:361:LYS:HB2	1:A:457:TYR:CD1	2.52	0.45
1:A:295[B]:LEU:CD2	1:A:387:PHE:CE2	3.00	0.44
1:A:27:TYR:O	1:A:247:VAL:HA	2.16	0.44
1:A:150:ASN:HD21	1:A:190:ASN:HB3	1.82	0.44
1:B:361:LYS:HB2	1:B:457:TYR:CD1	2.52	0.44
1:B:168:HIS:HB3	1:B:169:PRO:HA	1.99	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:296:HIS:HB3	1:B:365:PRO:HD2	1.99	0.44
1:A:361:LYS:HB2	1:A:457:TYR:CE1	2.53	0.43
1:B:249:SER:O	1:B:250:LYS:HB2	2.17	0.43
1:B:74:TYR:O	1:B:78[A]:GLN:HG2	2.18	0.43
1:B:31:ILE:HD13	1:B:51:VAL:CG1	2.48	0.43
1:B:361:LYS:HB2	1:B:457:TYR:CE1	2.53	0.43
1:B:260:ILE:HD12	1:B:385:VAL:HG11	2.00	0.43
1:A:249:SER:O	1:A:250:LYS:HB2	2.18	0.43
1:B:4:THR:HG21	1:B:191:LYS:HG3	1.99	0.43
1:B:237:ASP:OD2	1:B:243:HIS:CE1	2.67	0.43
1:A:195:HIS:HA	1:A:199:GLU:OE1	2.18	0.43
1:B:195:HIS:HA	1:B:199:GLU:OE1	2.18	0.43
1:B:300:ARG:HB2	1:B:457:TYR:HD1	1.83	0.42
1:A:338:PRO:HA	1:A:339:PRO:HD3	1.79	0.42
1:B:89:PHE:CZ	1:B:97:ASN:HB3	2.55	0.42
1:B:351:PRO:HB3	1:B:432:LEU:HD22	2.01	0.42
1:A:366:LEU:HD12	1:A:394:THR:HB	2.01	0.42
1:B:92[A]:GLU:OE2	1:B:384:ASN:HB3	2.19	0.42
1:A:374:LYS:HD2	10:A:547:HOH:O	2.19	0.42
1:A:147:TYR:CD1	1:A:147:TYR:N	2.88	0.42
1:A:162:PHE:CE2	1:A:409[A]:LEU:HD21	2.55	0.41
1:B:324:ASP:HB3	1:B:398:HIS:CG	2.55	0.41
1:A:326:TYR:CD1	1:A:496:SER:HA	2.56	0.41
1:A:110:ASN:HB2	3:A:523:FAD:N5	2.35	0.41
1:A:288:HIS:HB3	1:A:291:VAL:HG23	2.02	0.41
1:B:156:SER:O	1:B:160:THR:HG23	2.21	0.41
1:B:259:THR:OG1	1:B:461:HIS:HA	2.20	0.41
1:A:295[A]:LEU:HD22	1:A:463:GLY:HA2	2.03	0.40
1:A:273:PRO:HB2	1:A:276:TYR:HB3	2.03	0.40
1:A:324:ASP:HB3	1:A:398:HIS:CG	2.56	0.40
1:B:338:PRO:HA	1:B:339:PRO:HD3	1.89	0.40
1:A:32:VAL:HG13	1:A:217:VAL:HG21	2.04	0.40
1:B:380:ARG:NH2	10:B:1176:HOH:O	2.50	0.40
1:B:422:ASP:O	1:B:423:LEU:CD2	2.68	0.40
1:A:74:TYR:O	1:A:78[A]:GLN:HG2	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	529/521 (102%)	517 (98%)	11 (2%)	1 (0%)	56	28
1	B	527/521 (101%)	514 (98%)	13 (2%)	0	100	100
All	All	1056/1042 (101%)	1031 (98%)	24 (2%)	1 (0%)	59	31

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	486	ASP

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	453/443 (102%)	450 (99%)	3 (1%)	91	80
1	B	451/443 (102%)	447 (99%)	4 (1%)	87	73
All	All	904/886 (102%)	897 (99%)	7 (1%)	89	77

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	288	HIS
1	A	430	ASN
1	A	498	PRO
1	B	288	HIS
1	B	422	ASP
1	B	430	ASN
1	B	498	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	150	ASN
1	A	224	ASN
1	A	243	HIS
1	A	281	ASN
1	A	323	ASN
1	B	65	ASN
1	B	150	ASN
1	B	190	ASN
1	B	224	ASN
1	B	243	HIS
1	B	281	ASN
1	B	323	ASN
1	B	384	ASN
1	B	430	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

17 carbohydrates 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	$\# Z > 2$
5	NAG	A	525	1,5	12,14,15	0.53	0	15,19,21	0.81	0
5	NAG	A	526	5	12,14,15	0.60	0	15,19,21	0.85	0
5	FUL	A	527	5	9,10,11	0.41	0	10,14,16	0.30	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	MAN	A	528	5	10,11,12	0.48	0	11,15,17	0.58	0
5	MAN	A	529	5	10,11,12	0.46	0	11,15,17	0.25	0
6	NAG	A	530	1,6	12,14,15	0.49	0	15,19,21	0.72	0
6	NAG	A	531	6	12,14,15	0.40	0	15,19,21	0.74	0
6	FUC	A	532	6	9,10,11	0.42	0	10,14,16	0.25	0
7	NAG	B	524	7	12,14,15	0.45	0	15,19,21	0.72	0
7	NAG	B	525	1,7	12,14,15	0.51	0	15,19,21	0.70	0
7	FUC	B	526	7	9,10,11	0.42	0	10,14,16	0.27	0
7	BMA	B	527	7	10,11,12	0.39	0	11,15,17	0.27	0
7	MAN	B	528	7	10,11,12	0.39	0	11,15,17	0.27	0
8	NAG	B	529	1,8	12,14,15	0.50	0	15,19,21	1.00	0
8	NAG	B	530	8	12,14,15	0.55	0	15,19,21	0.70	0
8	BMA	B	531	8	10,11,12	0.42	0	11,15,17	0.47	0
8	MAN	B	532	8	10,11,12	0.40	0	11,15,17	0.24	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
5	NAG	A	525	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	526	5	-	0/6/23/26	0/1/1/1
5	FUL	A	527	5	-	0/0/17/20	0/1/1/1
5	MAN	A	528	5	-	0/2/19/22	0/1/1/1
5	MAN	A	529	5	-	0/2/19/22	0/1/1/1
6	NAG	A	530	1,6	-	0/6/23/26	0/1/1/1
6	NAG	A	531	6	-	0/6/23/26	0/1/1/1
6	FUC	A	532	6	-	0/0/17/20	0/1/1/1
7	NAG	B	524	7	-	0/6/23/26	0/1/1/1
7	NAG	B	525	1,7	-	0/6/23/26	0/1/1/1
7	FUC	B	526	7	-	0/0/17/20	0/1/1/1
7	BMA	B	527	7	-	0/2/19/22	0/1/1/1
7	MAN	B	528	7	-	0/2/19/22	0/1/1/1
8	NAG	B	529	1,8	-	0/6/23/26	0/1/1/1
8	NAG	B	530	8	-	0/6/23/26	0/1/1/1
8	BMA	B	531	8	-	0/2/19/22	0/1/1/1
8	MAN	B	532	8	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	IPA	A	522	-	3,3,3	0.52	0	3,3,3	0.87	0
3	FAD	A	523	-	58,58,58	2.17	19 (32%)	85,89,89	1.48	10 (11%)
4	NAG	A	524	1	12,14,15	0.46	0	15,19,21	0.65	0
4	NAG	A	533	1	12,14,15	0.41	0	15,19,21	0.74	0
2	IPA	B	522	-	3,3,3	0.51	0	3,3,3	0.88	0
3	FAD	B	523	-	58,58,58	2.21	19 (32%)	85,89,89	1.52	12 (14%)
9	NDG	B	533	-	12,14,15	0.34	0	15,19,21	0.97	1 (6%)
4	NAG	B	534	1	12,14,15	0.51	0	15,19,21	0.68	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	IPA	A	522	-	-	0/0/0/0	0/0/0/0
3	FAD	A	523	-	-	0/34/50/50	0/1/6/6
4	NAG	A	524	1	-	0/6/23/26	0/1/1/1
4	NAG	A	533	1	-	0/6/23/26	0/1/1/1
2	IPA	B	522	-	-	0/0/0/0	0/0/0/0
3	FAD	B	523	-	-	0/34/50/50	0/1/6/6
9	NDG	B	533	-	-	0/6/23/26	0/1/1/1
4	NAG	B	534	1	-	0/6/23/26	0/1/1/1

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	523	FAD	P-O3P	-6.82	1.47	1.59
3	A	523	FAD	P-O3P	-6.40	1.48	1.59
3	B	523	FAD	C9A-N10	5.65	1.47	1.38
3	A	523	FAD	C9A-N10	5.43	1.46	1.38
3	A	523	FAD	C1'-C2'	4.99	1.56	1.51
3	B	523	FAD	C1'-C2'	4.98	1.56	1.51
3	B	523	FAD	C4X-C10	4.35	1.48	1.40
3	A	523	FAD	C4X-C10	4.22	1.48	1.40
3	B	523	FAD	PA-O2A	-3.98	1.37	1.55
3	A	523	FAD	PA-O2A	-3.92	1.37	1.55
3	A	523	FAD	O4B-C1B	3.70	1.47	1.41
3	B	523	FAD	C4-C4X	3.54	1.47	1.41
3	B	523	FAD	O4B-C1B	3.42	1.46	1.41
3	B	523	FAD	P-O2P	-3.37	1.40	1.55
3	B	523	FAD	C2-N3	3.33	1.43	1.37
3	A	523	FAD	C4-C4X	3.27	1.46	1.41
3	A	523	FAD	P-O2P	-3.24	1.40	1.55
3	A	523	FAD	C2-N3	3.15	1.43	1.37
3	A	523	FAD	O5'-C5'	2.76	1.56	1.44
3	A	523	FAD	PA-O3P	2.67	1.64	1.59
3	B	523	FAD	O5'-C5'	2.66	1.55	1.44
3	B	523	FAD	C5X-N5	2.62	1.39	1.35
3	B	523	FAD	PA-O3P	2.56	1.64	1.59
3	A	523	FAD	C8-C7	2.53	1.48	1.40
3	A	523	FAD	C5X-N5	2.51	1.39	1.35
3	A	523	FAD	C2B-C1B	-2.47	1.49	1.53
3	B	523	FAD	C8-C7	2.44	1.47	1.40
3	B	523	FAD	C2B-C1B	-2.37	1.50	1.53
3	B	523	FAD	C4A-N3A	2.31	1.39	1.35
3	A	523	FAD	C4A-N3A	2.31	1.39	1.35
3	B	523	FAD	C2A-N3A	2.27	1.36	1.32
3	A	523	FAD	C2A-N1A	2.27	1.38	1.33
3	B	523	FAD	C6-C5X	2.21	1.44	1.41
3	A	523	FAD	C2A-N3A	2.19	1.36	1.32
3	B	523	FAD	O4B-C4B	2.16	1.50	1.45
3	A	523	FAD	O4B-C4B	2.16	1.50	1.45
3	B	523	FAD	C2A-N1A	2.09	1.38	1.33
3	A	523	FAD	C6-C5X	2.05	1.44	1.41

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	523	FAD	O4B-C1B-N9A	-7.16	101.78	108.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	523	FAD	O4B-C1B-N9A	-6.89	102.03	108.44
3	A	523	FAD	C2-N1-C10	5.66	120.69	114.98
3	B	523	FAD	C2-N1-C10	5.58	120.60	114.98
3	B	523	FAD	C4X-C10-N10	-3.97	118.53	120.51
3	A	523	FAD	C4X-C10-N10	-3.72	118.65	120.51
3	B	523	FAD	N3A-C2A-N1A	-3.41	125.86	128.71
3	A	523	FAD	N3A-C2A-N1A	-3.28	125.97	128.71
3	B	523	FAD	O5B-PA-O1A	-2.60	99.19	109.37
3	A	523	FAD	O5B-PA-O1A	-2.56	99.36	109.37
9	B	533	NDG	C3-C2-N2	-2.54	107.90	111.76
3	B	523	FAD	C5'-C4'-C3'	-2.36	107.61	112.06
3	A	523	FAD	C5'-C4'-C3'	-2.34	107.64	112.06
3	B	523	FAD	C5A-C4A-N9A	-2.32	103.81	107.16
3	B	523	FAD	C2A-N1A-C6A	2.30	122.93	118.77
3	B	523	FAD	C4A-C5A-N7A	2.29	111.48	109.52
3	A	523	FAD	C4A-C5A-N7A	2.27	111.47	109.52
3	A	523	FAD	C5A-C4A-N9A	-2.21	103.98	107.16
3	B	523	FAD	N3A-C4A-N9A	2.18	129.38	125.43
3	A	523	FAD	C2A-N1A-C6A	2.15	122.65	118.77
3	A	523	FAD	N3A-C4A-N9A	2.14	129.30	125.43
3	B	523	FAD	O4B-C1B-C2B	-2.10	103.55	106.77
3	B	523	FAD	C4B-O4B-C1B	-2.03	107.55	109.75

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	521/521 (100%)	-0.22	5 (0%) 79 82	6, 11, 21, 36	2 (0%)
1	B	521/521 (100%)	-0.15	5 (0%) 79 82	7, 12, 24, 40	3 (0%)
All	All	1042/1042 (100%)	-0.19	10 (0%) 79 82	6, 11, 23, 40	5 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	124	ALA	3.3
1	B	422	ASP	2.8
1	B	423	LEU	2.4
1	A	521	ASP	2.4
1	B	23	LEU	2.4
1	B	424	PRO	2.4
1	A	347	SER	2.4
1	B	425	GLY	2.3
1	A	519	ALA	2.1
1	A	520	SER	2.0

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

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

6.3 Carbohydrates ⓘ

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	MAN	A	529	11/12	0.23	40.09	52,52,53,53	0
8	MAN	B	532	11/12	0.26	26.80	50,50,52,52	0
5	NAG	A	526	14/15	0.21	10.64	30,33,39,43	0
8	NAG	B	530	14/15	0.16	4.78	27,30,35,38	0
6	NAG	A	530	14/15	0.13	4.58	24,29,35,36	0
7	NAG	B	525	14/15	0.10	2.47	18,22,25,27	0
7	MAN	B	528	11/12	0.24	1.43	43,45,46,46	0
8	NAG	B	529	14/15	0.10	1.30	14,19,22,23	0
5	NAG	A	525	14/15	0.10	0.91	14,20,25,29	0
7	BMA	B	527	11/12	0.20	-	38,40,41,42	0
5	MAN	A	528	11/12	0.29	-	46,48,50,52	0
6	NAG	A	531	14/15	0.29	-	40,42,43,44	0
7	FUC	B	526	10/11	0.25	-	31,34,35,36	0
6	FUC	A	532	10/11	0.31	-	39,41,42,43	0
5	FUL	A	527	10/11	0.39	-	35,39,41,41	0
7	NAG	B	524	14/15	0.17	-	28,30,35,36	0
8	BMA	B	531	11/12	0.23	-	42,44,47,49	0

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	NAG	A	533	14/15	0.27	16.18	37,41,42,43	0
9	NDG	B	533	14/15	0.26	4.41	40,43,45,45	0
2	IPA	A	522	4/4	0.11	3.98	19,20,20,21	0
4	NAG	A	524	14/15	0.12	2.20	15,17,21,22	0
4	NAG	B	534	14/15	0.11	1.33	15,21,24,26	0
2	IPA	B	522	4/4	0.08	0.88	19,20,21,21	0
3	FAD	B	523	53/53	0.07	-0.22	5,7,8,9	0
3	FAD	A	523	53/53	0.07	-0.25	4,6,9,9	0

6.5 Other polymers ⓘ

There are no such residues in this entry.