Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 07:46 PM GMT

PDB ID	:	1H6B
Title	:	REDUCED PRECURSOR FORM OF GLUCOSE-FRUCTOSE OXIDOREDUCTASE FROM ZYMOMONAS MOBILIS COMPLEXED WITH GLYCEROL
Authors	:	Nurizzo, D.; Baker, E.N.
Deposited on	:	2001-06-11
Resolution	:	2.60 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:

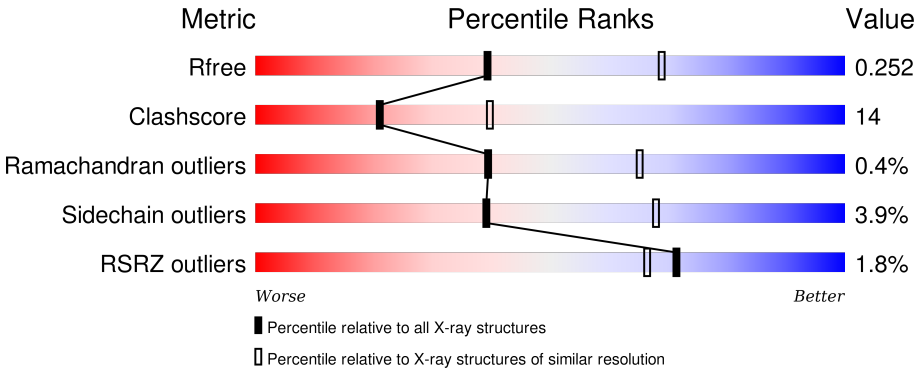
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance

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

The reported resolution of this entry is 2.60 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	433	66% 20% • 12%
1	B	433	3% 61% 24% • 12%

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	600	-	-	-	X

2 Entry composition

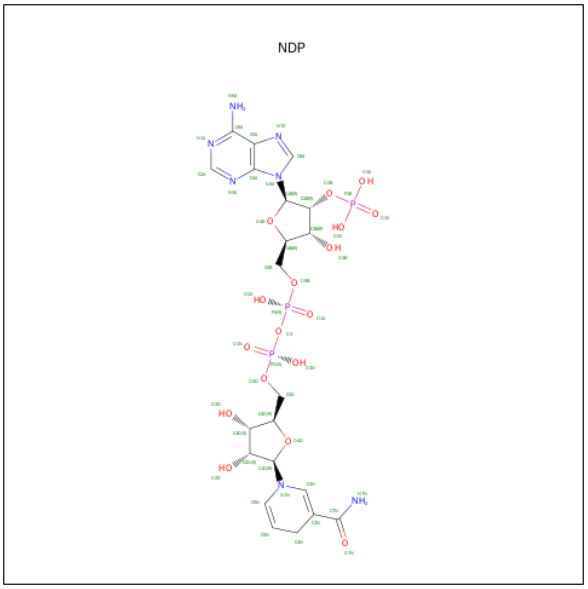
There are 4 unique types of molecules in this entry. The entry contains 6324 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PRECURSOR FORM OF GLUCOSE-FRUCTOSE OXIDOREDUCTASE.

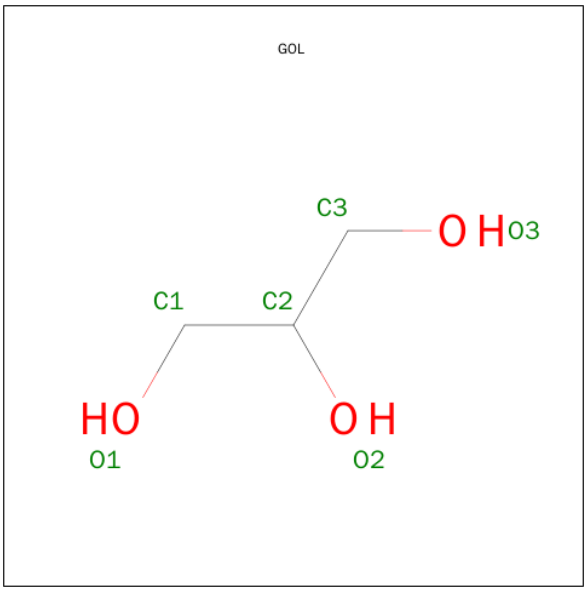
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	381	Total	C	N	O	S	0	0	0
			2960	1856	530	554	20			
1	B	381	Total	C	N	O	S	0	0	0
			2960	1856	530	554	20			

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0

- Molecule 4 is water.

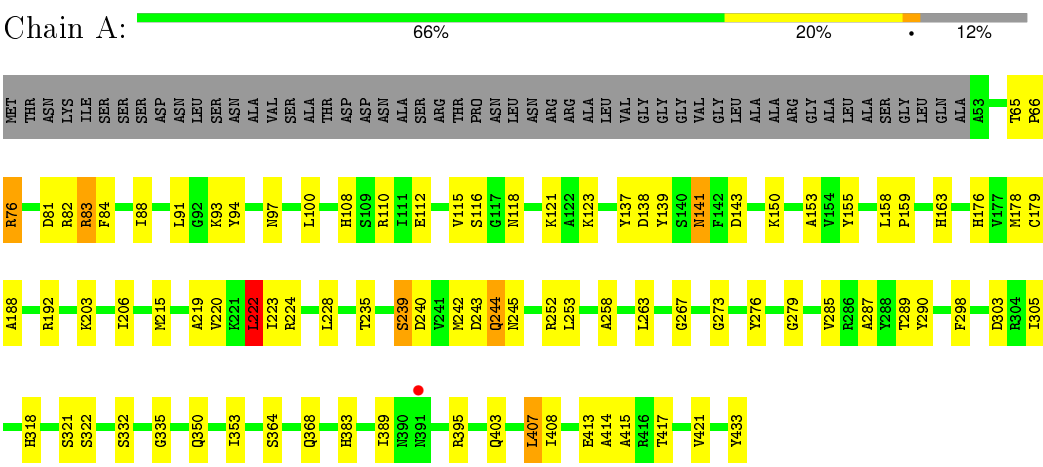
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	203	Total O 203 203	0	0
4	B	93	Total O 93 93	0	0

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.

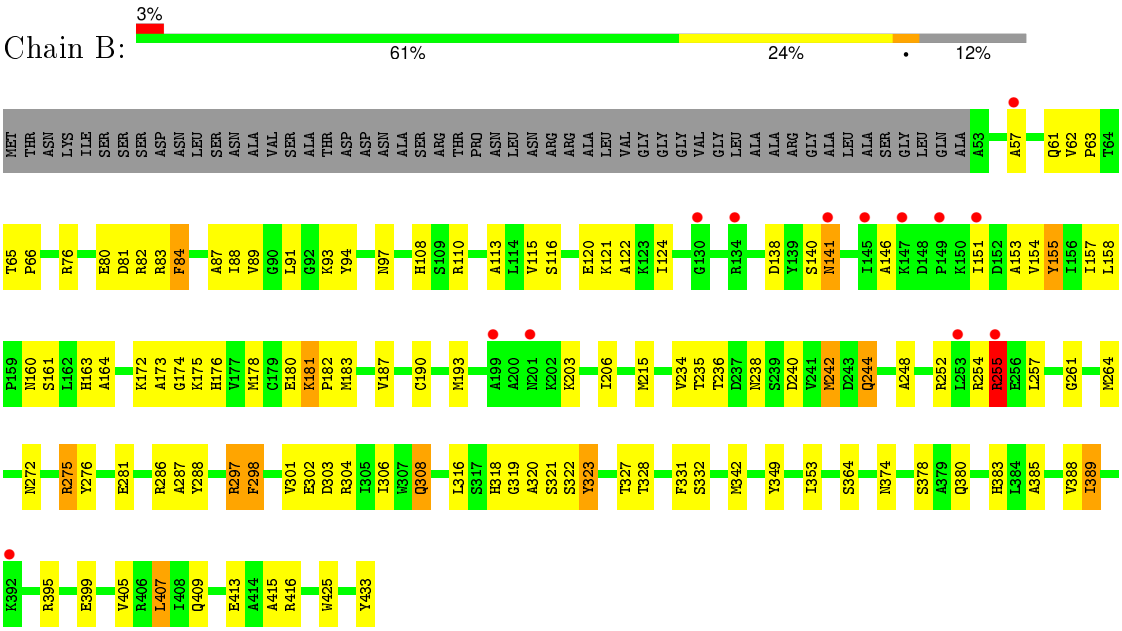
- Molecule 1: PRECURSOR FORM OF GLUCOSE-FRUCTOSE OXIDOREDUCTASE

Chain A:



- Molecule 1: PRECURSOR FORM OF GLUCOSE-FRUCTOSE OXIDOREDUCTASE

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	83.80Å 92.77Å 115.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.60 14.99 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.3 (15.00-2.60) 97.4 (14.99-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.81 (at 2.61Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.200 , 0.252 0.200 , 0.252	Depositor DCC
R_{free} test set	806 reflections (2.94%)	DCC
Wilson B-factor (Å^2)	38.1	Xtriage
Anisotropy	0.495	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 55.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 27442 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6324	wwPDB-VP
Average B, all atoms (Å^2)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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: GOL, NDP

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	0.38	0/3024	0.64	2/4093 (0.0%)
1	B	0.48	0/3024	0.73	8/4093 (0.2%)
All	All	0.43	0/6048	0.69	10/8186 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	286	ARG	NE-CZ-NH2	6.78	123.69	120.30
1	B	254	ARG	NE-CZ-NH2	6.76	123.68	120.30
1	B	252	ARG	NE-CZ-NH2	6.75	123.67	120.30
1	B	255	ARG	NE-CZ-NH2	6.73	123.67	120.30
1	B	275	ARG	NE-CZ-NH2	6.34	123.47	120.30
1	A	222	LEU	CA-CB-CG	6.09	129.31	115.30
1	B	297	ARG	NE-CZ-NH2	5.94	123.27	120.30
1	B	264	MET	CG-SD-CE	5.80	109.48	100.20
1	B	242	MET	CG-SD-CE	5.50	109.00	100.20
1	A	178	MET	CG-SD-CE	5.06	108.30	100.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	298	PHE	Mainchain

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	2960	0	2920	76	0
1	B	2960	0	2920	100	0
2	A	48	0	26	2	0
2	B	48	0	26	5	0
3	A	6	0	8	0	0
3	B	6	0	8	0	0
4	A	203	0	0	2	0
4	B	93	0	0	1	0
All	All	6324	0	5908	171	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 14.

All (171) 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:A:138:ASP:H	1:A:141:ASN:HD21	1.21	0.89
1:A:244:GLN:HE21	1:A:244:GLN:H	1.26	0.84
1:B:87:ALA:HB3	1:B:154:VAL:HG12	1.59	0.83
1:A:176:HIS:HD2	1:A:203:LYS:H	1.29	0.81
1:B:255:ARG:HB2	1:B:255:ARG:NH1	1.94	0.81
1:B:176:HIS:HD2	1:B:203:LYS:H	1.26	0.80
1:B:153:ALA:HB2	1:B:176:HIS:HB2	1.63	0.80
1:A:219:ALA:O	1:A:222:LEU:HD22	1.84	0.78
1:A:244:GLN:NE2	1:A:244:GLN:H	1.82	0.78
1:A:306:ILE:HD13	1:B:318:HIS:HB3	1.66	0.75
1:A:383:HIS:HE1	1:A:395:ARG:H	1.37	0.73
1:A:69:ARG:HB3	1:A:70:PRO:HD2	1.71	0.72
1:B:374:ASN:ND2	1:B:378:SER:HB2	2.06	0.70
1:B:255:ARG:HB2	1:B:255:ARG:HH11	1.55	0.70
1:A:222:LEU:HD23	1:A:223:ILE:N	2.06	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:383:HIS:HE1	1:B:395:ARG:H	1.42	0.68
1:B:57:ALA:HB1	1:B:61:GLN:HE22	1.58	0.68
1:B:57:ALA:HB1	1:B:61:GLN:NE2	2.09	0.68
1:B:84:PHE:HE1	1:B:388:VAL:HG21	1.58	0.67
1:B:383:HIS:CE1	1:B:395:ARG:H	2.12	0.67
1:B:176:HIS:CD2	1:B:203:LYS:H	2.11	0.67
1:A:383:HIS:CE1	1:A:395:ARG:H	2.13	0.67
1:B:89:VAL:HA	1:B:115:VAL:HB	1.77	0.67
1:A:83:ARG:HD2	1:A:389:ILE:HD11	1.76	0.67
1:B:181:LYS:HD2	1:B:181:LYS:C	2.15	0.67
1:B:138:ASP:H	1:B:141:ASN:HD21	1.41	0.66
1:B:255:ARG:CB	1:B:255:ARG:HH11	2.08	0.66
1:B:181:LYS:O	1:B:181:LYS:HD2	1.96	0.65
1:A:115:VAL:HG22	1:A:137:TYR:HB2	1.78	0.64
1:B:116:SER:HB3	1:B:122:ALA:HB2	1.80	0.64
1:A:138:ASP:N	1:A:141:ASN:HD21	1.95	0.60
1:B:413:GLU:HG3	1:B:416:ARG:HH21	1.66	0.59
1:B:244:GLN:HE21	1:B:244:GLN:H	1.49	0.58
1:B:388:VAL:HG23	1:B:389:ILE:N	2.18	0.58
1:B:89:VAL:HG12	1:B:158:LEU:HD21	1.86	0.58
1:A:215:MET:HE1	1:A:364:SER:HB3	1.86	0.57
1:A:239:SER:HB3	1:A:322:SER:HB3	1.87	0.57
1:A:242:MET:SD	1:A:253:LEU:HD21	2.45	0.57
1:B:238:ASN:O	1:B:321:SER:HA	2.06	0.55
1:B:261:GLY:CA	1:B:301:VAL:HB	2.36	0.55
1:B:304:ARG:HG2	1:B:322:SER:HB2	1.88	0.55
1:B:389:ILE:O	1:B:389:ILE:HG22	2.06	0.55
1:B:238:ASN:HA	1:B:328:THR:O	2.06	0.55
1:A:153:ALA:HB2	1:A:176:HIS:HB2	1.89	0.55
1:A:318:HIS:HB3	1:B:306:ILE:HD13	1.89	0.54
1:B:331:PHE:HB2	1:B:342:MET:HB2	1.90	0.54
1:A:83:ARG:HD2	1:A:389:ILE:CD1	2.38	0.54
1:A:308:GLN:OE1	1:B:288:TYR:HB3	2.08	0.53
1:B:81:ASP:CG	1:B:83:ARG:HG3	2.28	0.53
1:A:223:ILE:HG12	1:A:228:LEU:HD22	1.91	0.53
1:B:84:PHE:CE1	1:B:388:VAL:HG21	2.41	0.53
1:A:279:GLY:HA2	1:A:433:TYR:CD2	2.43	0.53
1:A:413:GLU:O	1:A:417:THR:HG23	2.09	0.53
1:A:285:VAL:HG12	1:A:309:MET:HG2	1.91	0.53
1:B:121:LYS:HE3	2:B:500:NDP:O2X	2.09	0.52
1:A:289:THR:HB	1:A:305:ILE:HG13	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:LEU:CD2	1:A:223:ILE:HG13	2.40	0.52
1:B:138:ASP:N	1:B:141:ASN:HD21	2.06	0.52
1:B:146:ALA:HB2	1:B:173:ALA:HA	1.91	0.52
1:B:244:GLN:HE22	1:B:297:ARG:HD3	1.76	0.51
1:B:158:LEU:O	1:B:163:HIS:NE2	2.44	0.51
1:A:81:ASP:OD1	1:A:108:HIS:HA	2.12	0.50
1:B:244:GLN:NE2	1:B:297:ARG:HD3	2.27	0.49
1:B:235:THR:HB	1:B:332:SER:HB3	1.94	0.49
1:A:181:LYS:C	1:A:181:LYS:HD2	2.32	0.49
1:A:94:TYR:CD1	2:A:500:NDP:H41N	2.47	0.49
1:B:120:GLU:O	1:B:124:ILE:HG13	2.12	0.49
1:B:388:VAL:HG23	1:B:389:ILE:H	1.78	0.49
1:B:80:GLU:OE1	1:B:108:HIS:HE1	1.94	0.49
1:B:407:LEU:HD13	1:B:425:TRP:CZ2	2.48	0.49
1:A:235:THR:HB	1:A:332:SER:HB3	1.95	0.49
1:A:350:GLN:HB3	4:A:2147:HOH:O	2.12	0.49
1:B:183:MET:HE3	1:B:206:ILE:HD11	1.94	0.49
1:B:413:GLU:HG3	1:B:416:ARG:NH2	2.28	0.49
1:B:63:PRO:C	1:B:65:THR:H	2.16	0.49
1:B:173:ALA:O	1:B:175:LYS:HE3	2.13	0.48
1:A:110:ARG:HH21	1:A:112:GLU:CG	2.26	0.48
1:A:306:ILE:HD12	1:B:308:GLN:HG2	1.96	0.48
1:B:82:ARG:HB3	1:B:110:ARG:HB3	1.95	0.48
1:A:244:GLN:HE21	1:A:244:GLN:N	2.03	0.48
1:B:138:ASP:H	1:B:141:ASN:ND2	2.09	0.48
1:A:93:LYS:O	1:A:97:ASN:HB2	2.14	0.48
1:B:255:ARG:CB	1:B:255:ARG:NH1	2.70	0.47
1:A:276:TYR:CB	1:A:433:TYR:HB3	2.44	0.47
1:A:179:CYS:O	1:A:206:ILE:HA	2.15	0.47
1:B:374:ASN:HD21	1:B:378:SER:HB2	1.77	0.47
1:B:255:ARG:HD3	4:B:2041:HOH:O	2.14	0.47
1:A:267:GLY:CA	1:A:321:SER:HB3	2.44	0.47
1:A:116:SER:O	1:A:138:ASP:HA	2.15	0.47
1:B:116:SER:CB	1:B:122:ALA:HB2	2.44	0.47
1:B:174:GLY:O	1:B:175:LYS:HD3	2.16	0.46
1:B:405:VAL:O	1:B:409:GLN:HG2	2.14	0.46
1:B:287:ALA:O	1:B:415:ALA:HA	2.15	0.46
1:A:222:LEU:HD23	1:A:223:ILE:HG13	1.98	0.46
1:B:302:GLU:HG2	1:B:302:GLU:O	2.16	0.46
1:B:84:PHE:O	1:B:110:ARG:HG2	2.15	0.46
1:A:83:ARG:HG3	1:A:84:PHE:N	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:ASN:HD22	1:A:141:ASN:C	2.18	0.45
1:A:76:ARG:HG2	1:A:76:ARG:HH11	1.80	0.45
1:B:141:ASN:H	1:B:141:ASN:ND2	2.14	0.45
1:A:239:SER:CB	1:A:322:SER:HB3	2.47	0.45
1:B:236:THR:O	1:B:319:GLY:HA2	2.16	0.45
1:A:158:LEU:O	1:A:163:HIS:NE2	2.47	0.45
1:A:287:ALA:O	1:A:415:ALA:HA	2.17	0.45
1:A:88:ILE:HG13	1:A:100:LEU:HD21	1.98	0.45
1:A:69:ARG:HB3	1:A:70:PRO:CD	2.45	0.45
1:B:160:ASN:OD1	1:B:182:PRO:HD2	2.17	0.45
1:A:181:LYS:HA	1:A:182:PRO:C	2.36	0.45
1:B:298:PHE:CG	1:B:303:ASP:HB2	2.52	0.44
1:A:176:HIS:CD2	1:A:203:LYS:H	2.20	0.44
1:B:234:VAL:HA	1:B:332:SER:O	2.17	0.44
1:A:273:GLY:HA2	1:A:276:TYR:CE2	2.52	0.44
1:B:62:VAL:HA	1:B:63:PRO:HD3	1.83	0.44
1:B:93:LYS:O	1:B:97:ASN:HB2	2.18	0.44
1:B:374:ASN:HD21	1:B:378:SER:CB	2.31	0.44
1:B:138:ASP:OD1	1:B:140:SER:N	2.40	0.44
1:B:172:LYS:C	1:B:174:GLY:H	2.20	0.44
1:A:395:ARG:HG3	1:A:395:ARG:O	2.19	0.43
1:A:158:LEU:HB3	1:A:159:PRO:HD2	2.00	0.43
1:A:118:ASN:HB3	1:A:121:LYS:HE2	1.99	0.43
1:A:298:PHE:CD2	1:A:303:ASP:HB2	2.54	0.43
1:A:65:THR:HB	1:A:66:PRO:HD2	2.00	0.43
1:B:153:ALA:CB	1:B:176:HIS:HB2	2.43	0.43
1:A:228:LEU:O	1:A:335:GLY:HA3	2.18	0.43
1:B:248:ALA:HA	2:B:500:NDP:O1A	2.19	0.43
1:B:178:MET:HE3	1:B:380:GLN:HG2	2.00	0.43
1:B:272:ASN:O	1:B:275:ARG:HB2	2.18	0.43
1:B:187:VAL:O	1:B:190:CYS:HB2	2.19	0.43
1:B:91:LEU:HD12	1:B:121:LYS:HD2	2.00	0.43
1:B:298:PHE:CD2	1:B:303:ASP:HB2	2.54	0.43
1:B:276:TYR:CB	1:B:433:TYR:HB3	2.49	0.43
1:B:164:ALA:HA	1:B:193:MET:SD	2.58	0.43
1:B:261:GLY:HA3	1:B:301:VAL:HB	2.01	0.42
1:B:215:MET:HE1	1:B:364:SER:HB3	2.01	0.42
1:A:290:TYR:HB3	1:B:316:LEU:HD11	2.01	0.42
1:A:123:LYS:HB2	1:A:123:LYS:HE3	1.84	0.42
1:B:88:ILE:HD12	1:B:155:TYR:HB3	2.02	0.42
1:B:399:GLU:OE1	1:B:399:GLU:N	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:353:ILE:CG1	1:A:364:SER:HB2	2.49	0.42
1:B:161:SER:OG	1:B:257:LEU:HD23	2.20	0.42
1:A:82:ARG:HB3	1:A:110:ARG:HB3	2.01	0.42
1:A:112:GLU:HG3	1:A:150:LYS:O	2.19	0.42
1:B:180:GLU:OE1	2:B:500:NDP:H2N	2.19	0.42
1:A:110:ARG:NH2	1:A:112:GLU:CG	2.83	0.42
1:B:161:SER:HB3	1:B:257:LEU:HA	2.02	0.42
1:B:157:ILE:HG22	2:B:500:NDP:H51N	2.01	0.41
1:A:243:ASP:OD1	1:A:245:ASN:HB2	2.20	0.41
1:A:252:ARG:HA	1:A:258:ALA:HB2	2.01	0.41
1:A:220:VAL:O	1:A:224:ARG:HG2	2.20	0.41
1:B:151:ILE:O	1:B:175:LYS:HD2	2.20	0.41
1:B:141:ASN:N	1:B:141:ASN:ND2	2.69	0.41
1:A:298:PHE:CG	1:A:303:ASP:HB2	2.54	0.41
1:B:242:MET:HE1	1:B:323:TYR:CE1	2.56	0.41
1:A:383:HIS:HD2	4:A:2075:HOH:O	2.02	0.41
1:B:94:TYR:CD1	2:B:500:NDP:H41N	2.55	0.41
1:A:188:ALA:O	1:A:192:ARG:HG3	2.21	0.41
1:B:306:ILE:HG12	1:B:320:ALA:HB2	2.02	0.41
1:B:353:ILE:CG1	1:B:364:SER:HB2	2.51	0.41
1:A:403:GLN:O	1:A:407:LEU:HB2	2.21	0.41
1:A:83:ARG:CD	1:A:389:ILE:HD11	2.47	0.41
1:A:115:VAL:HA	1:A:137:TYR:O	2.20	0.41
1:A:91:LEU:HG	1:A:116:SER:HB2	2.01	0.40
1:B:385:ALA:O	1:B:389:ILE:HG13	2.21	0.40
1:B:308:GLN:O	1:B:308:GLN:NE2	2.55	0.40
1:A:139:TYR:OH	2:A:500:NDP:H1B	2.21	0.40
1:A:414:ALA:HB2	1:A:421:VAL:HG23	2.03	0.40
1:A:263:LEU:HD23	1:A:408:ILE:HG23	2.04	0.40
1:B:87:ALA:HA	1:B:113:ALA:O	2.22	0.40
1:B:181:LYS:HA	1:B:182:PRO:C	2.41	0.40
1:B:141:ASN:HD22	1:B:141:ASN:N	2.19	0.40
1:B:275:ARG:NE	1:B:281:GLU:OE1	2.42	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	379/433 (88%)	357 (94%)	22 (6%)	0	100 100
1	B	379/433 (88%)	350 (92%)	26 (7%)	3 (1%)	24 46
All	All	758/866 (88%)	707 (93%)	48 (6%)	3 (0%)	39 65

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	327	THR
1	B	389	ILE
1	B	66	PRO

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	308/345 (89%)	296 (96%)	12 (4%)	39 68
1	B	308/345 (89%)	296 (96%)	12 (4%)	39 68
All	All	616/690 (89%)	592 (96%)	24 (4%)	39 68

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	ARG
1	A	83	ARG
1	A	141	ASN

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Mol	Chain	Res	Type
1	A	143	ASP
1	A	155	TYR
1	A	181	LYS
1	A	222	LEU
1	A	239	SER
1	A	240	ASP
1	A	244	GLN
1	A	368	GLN
1	A	407	LEU
1	B	76	ARG
1	B	84	PHE
1	B	141	ASN
1	B	155	TYR
1	B	181	LYS
1	B	240	ASP
1	B	244	GLN
1	B	255	ARG
1	B	308	GLN
1	B	323	TYR
1	B	349	TYR
1	B	407	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	141	ASN
1	A	176	HIS
1	A	201	ASN
1	A	216	ASN
1	A	227	GLN
1	A	244	GLN
1	A	245	ASN
1	A	250	GLN
1	A	360	HIS
1	A	362	ASN
1	A	383	HIS
1	A	390	ASN
1	B	61	GLN
1	B	98	GLN
1	B	107	GLN
1	B	108	HIS
1	B	141	ASN

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Mol	Chain	Res	Type
1	B	176	HIS
1	B	216	ASN
1	B	227	GLN
1	B	238	ASN
1	B	244	GLN
1	B	245	ASN
1	B	250	GLN
1	B	308	GLN
1	B	362	ASN
1	B	374	ASN
1	B	375	ASN
1	B	383	HIS

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	NDP	A	500	-	42,52,52	1.63	8 (19%)	55,80,80	2.08	10 (18%)
3	GOL	A	600	-	5,5,5	0.96	0	5,5,5	0.42	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	NDP	B	500	-	42,52,52	1.62	10 (23%)	55,80,80	1.99	10 (18%)
3	GOL	B	600	-	5,5,5	0.93	0	5,5,5	0.40	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	NDP	A	500	-	-	0/30/77/77	0/5/5/5
3	GOL	A	600	-	-	0/4/4/4	0/0/0/0
2	NDP	B	500	-	-	0/30/77/77	0/5/5/5
3	GOL	B	600	-	-	0/4/4/4	0/0/0/0

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	NDP	P2B-O2X	-4.99	1.36	1.54
2	B	500	NDP	P2B-O2X	-4.91	1.37	1.54
2	A	500	NDP	C4N-C5N	-2.20	1.44	1.49
2	B	500	NDP	C4N-C5N	-2.08	1.44	1.49
2	B	500	NDP	C4A-N3A	2.01	1.38	1.35
2	B	500	NDP	O4D-C1D	2.06	1.47	1.42
2	A	500	NDP	O4D-C1D	2.08	1.47	1.42
2	A	500	NDP	P2B-O1X	2.12	1.58	1.51
2	B	500	NDP	P2B-O1X	2.20	1.58	1.51
2	B	500	NDP	C6N-N1N	2.34	1.44	1.37
2	A	500	NDP	C6N-N1N	2.54	1.45	1.37
2	B	500	NDP	P2B-O2B	2.69	1.68	1.60
2	B	500	NDP	O4B-C1B	2.74	1.44	1.41
2	B	500	NDP	C2N-C3N	2.80	1.41	1.34
2	A	500	NDP	C2N-C3N	2.84	1.41	1.34
2	A	500	NDP	O4B-C1B	3.20	1.45	1.41
2	B	500	NDP	C6N-C5N	4.02	1.41	1.33
2	A	500	NDP	C6N-C5N	4.42	1.41	1.33

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	500	NDP	N3A-C2A-N1A	-10.59	120.79	128.89
2	B	500	NDP	N3A-C2A-N1A	-10.38	120.94	128.89
2	B	500	NDP	C3N-C2N-N1N	-3.40	118.28	123.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	500	NDP	C3N-C2N-N1N	-3.18	118.58	123.14
2	A	500	NDP	C4N-C5N-C6N	-3.08	117.50	122.58
2	B	500	NDP	C4N-C5N-C6N	-2.97	117.68	122.58
2	B	500	NDP	C1D-N1N-C2N	-2.44	116.67	120.91
2	A	500	NDP	C1D-N1N-C2N	-2.13	117.19	120.91
2	B	500	NDP	O4B-C1B-N9A	2.01	112.30	108.10
2	B	500	NDP	O2X-P2B-O1X	2.01	117.05	110.58
2	A	500	NDP	O2X-P2B-O1X	2.04	117.14	110.58
2	A	500	NDP	PN-O3-PA	2.08	138.57	132.73
2	B	500	NDP	C4A-C5A-N7A	2.33	111.62	109.48
2	B	500	NDP	P2B-O2B-C2B	2.49	127.54	121.56
2	A	500	NDP	C4A-C5A-N7A	2.67	111.94	109.48
2	A	500	NDP	C5N-C4N-C3N	3.36	121.79	112.52
2	B	500	NDP	C5N-C4N-C3N	3.39	121.86	112.52
2	A	500	NDP	O4B-C1B-N9A	3.62	115.68	108.10
2	B	500	NDP	O3-PA-O5B	4.17	114.01	102.94
2	A	500	NDP	O3-PA-O5B	5.20	116.74	102.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	NDP	2	0
2	B	500	NDP	5	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	381/433 (87%)	-0.59	1 (0%) 94 93	19, 29, 41, 49	0
1	B	381/433 (87%)	-0.08	13 (3%) 49 41	23, 45, 64, 75	0
All	All	762/866 (87%)	-0.33	14 (1%) 71 66	19, 34, 60, 75	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	134	ARG	4.5
1	A	391	ASN	3.0
1	B	130	GLY	2.8
1	B	149	PRO	2.6
1	B	147	LYS	2.6
1	B	151	ILE	2.5
1	B	57	ALA	2.5
1	B	145	ILE	2.5
1	B	201	ASN	2.4
1	B	141	ASN	2.3
1	B	255	ARG	2.1
1	B	392	LYS	2.1
1	B	253	LEU	2.0
1	B	199	ALA	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

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
3	GOL	A	600	6/6	0.91	0.17	2.78	54,54,54,57	0
3	GOL	B	600	6/6	0.91	0.16	0.25	66,69,69,70	0
2	NDP	A	500	48/48	0.97	0.13	0.09	24,29,31,32	0
2	NDP	B	500	48/48	0.95	0.15	-0.32	44,49,51,53	0

6.5 Other polymers (i)

There are no such residues in this entry.