



# Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 11:49 pm BST

PDB ID : 1H6A  
Title : Reduced Precursor Form of Glucose-Fructose Oxidoreductase from *Zymomonas mobilis*  
Authors : Nurizzo, D.; Baker, E.N.  
Deposited on : 2001-06-11  
Resolution : 2.50 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

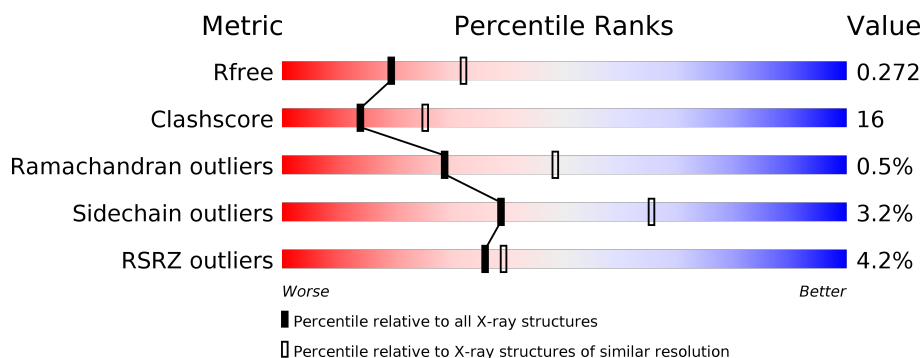
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. 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	433	
1	B	433	

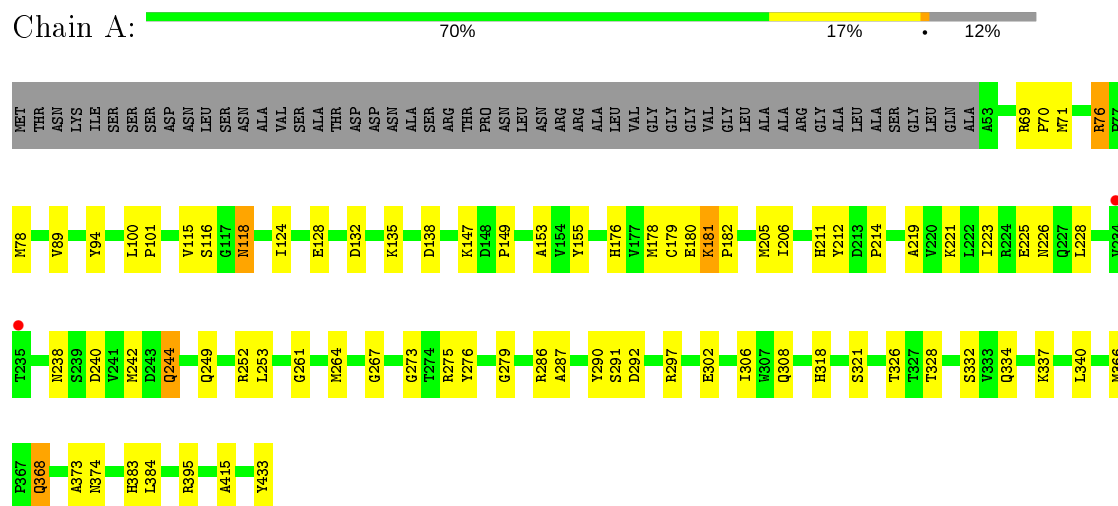


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	213	Total 213	O 213	0	0
3	B	85	Total 85	O 85	0	0

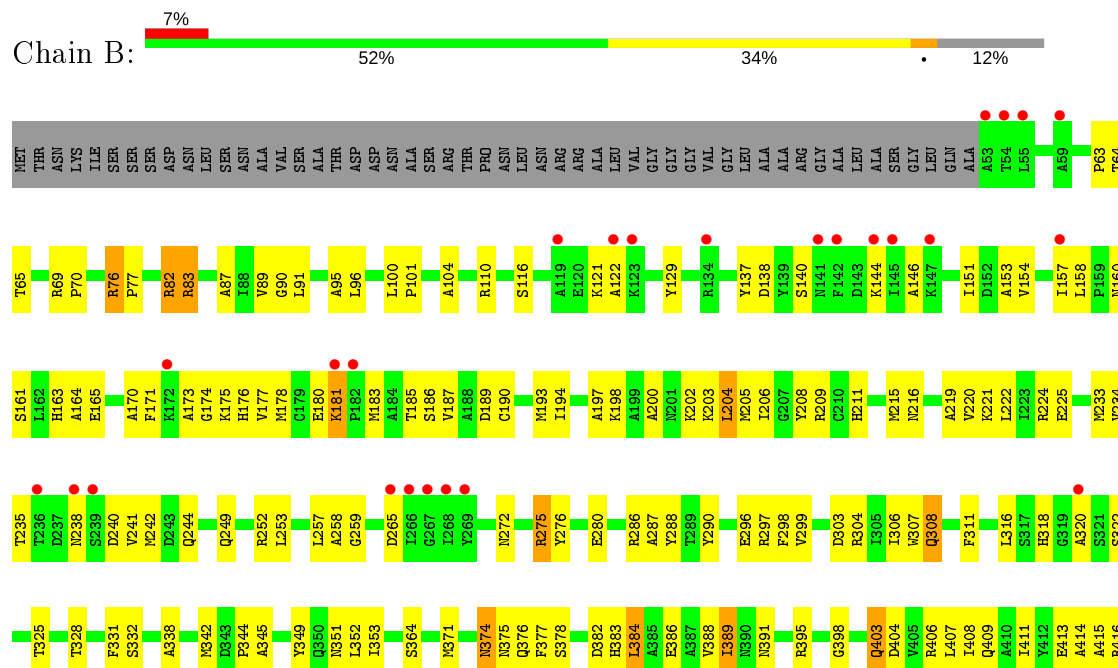
### 3 Residue-property plots

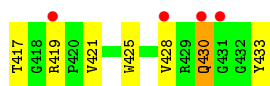
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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



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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 <sub>1</sub> 2 <sub>1</sub> 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.90 Å 93.06 Å 115.49 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.50 33.26 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.5 (15.00-2.50) 97.6 (33.26-2.50)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.07 (at 2.51 Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.204 , 0.275 0.197 , 0.272	Depositor DCC
$R_{free}$ test set	936 reflections (3.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	44.4	Xtriage
Anisotropy	0.650	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 46.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6314	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 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	0.41	0/3024	0.67	2/4093 (0.0%)
1	B	0.38	0/3024	0.64	3/4093 (0.1%)
All	All	0.39	0/6048	0.65	5/8186 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	82	ARG	NE-CZ-NH2	7.79	124.20	120.30
1	B	83	ARG	NE-CZ-NH2	7.34	123.97	120.30
1	A	275	ARG	NE-CZ-NH2	6.61	123.61	120.30
1	B	275	ARG	NE-CZ-NH2	5.25	122.93	120.30
1	A	71	MET	CG-SD-CE	5.17	108.47	100.20

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	2960	0	2920	60	0
1	B	2960	0	2920	136	0
2	A	48	0	26	4	0
2	B	48	0	26	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	213	0	0	3	0
3	B	85	0	0	2	0
All	All	6314	0	5892	188	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 16.

All (188) 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:368:GLN:NE2	1:A:368:GLN:H	1.65	0.94
1:A:205:MET:CE	1:A:384:LEU:HB2	1.99	0.93
1:A:244:GLN:NE2	1:A:244:GLN:H	1.66	0.92
1:A:368:GLN:N	1:A:368:GLN:HE21	1.69	0.90
1:A:244:GLN:HE21	1:A:244:GLN:N	1.70	0.89
1:B:153:ALA:HB2	1:B:176:HIS:HB2	1.55	0.89
1:B:211:HIS:HE1	1:B:272:ASN:HD21	1.22	0.88
1:A:244:GLN:HE21	1:A:244:GLN:H	0.87	0.86
1:B:413:GLU:HG3	1:B:416:ARG:HH12	1.46	0.80
1:A:306:ILE:HD13	1:B:318:HIS:HB3	1.65	0.78
1:B:76:ARG:HG2	1:B:77:PRO:HD2	1.65	0.77
1:B:211:HIS:CE1	1:B:272:ASN:HD21	2.02	0.77
1:A:368:GLN:HE21	1:A:368:GLN:H	0.85	0.76
1:B:89:VAL:HG12	1:B:158:LEU:HD21	1.66	0.76
1:B:176:HIS:HD2	1:B:203:LYS:HB2	1.51	0.76
1:B:403:GLN:HA	1:B:403:GLN:HE21	1.50	0.75
1:B:82:ARG:HG3	1:B:110:ARG:HB3	1.67	0.75
1:A:181:LYS:HD2	1:A:181:LYS:O	1.87	0.75
1:A:205:MET:HE1	1:A:384:LEU:HB2	1.70	0.73
1:B:209:ARG:HH11	1:B:376:GLN:HE22	1.37	0.72
1:B:221:LYS:O	1:B:225:GLU:HG3	1.89	0.71
1:A:118:ASN:HD22	1:A:118:ASN:C	1.94	0.70
1:A:181:LYS:C	1:A:181:LYS:HD2	2.15	0.67
1:A:242:MET:SD	1:A:249:GLN:HB3	2.33	0.67
1:B:221:LYS:HE2	1:B:225:GLU:OE2	1.95	0.66
1:B:76:ARG:HG2	1:B:77:PRO:CD	2.24	0.66
1:B:280:GLU:HG3	1:B:311:PHE:HD2	1.61	0.66
1:A:205:MET:HE2	1:A:384:LEU:HB2	1.79	0.65
1:B:204:LEU:HD13	1:B:398:GLY:HA3	1.78	0.65
1:B:63:PRO:C	1:B:65:THR:H	2.00	0.64
1:B:194:ILE:HG22	1:B:198:LYS:HE2	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:LEU:HB2	1:A:101:PRO:HD3	1.78	0.63
1:B:209:ARG:HH11	1:B:376:GLN:NE2	1.95	0.63
1:B:407:LEU:O	1:B:411:ILE:HG13	1.98	0.63
1:A:153:ALA:HB2	1:A:176:HIS:HB2	1.81	0.63
1:B:151:ILE:O	1:B:175:LYS:HD2	1.99	0.63
1:B:304:ARG:HG2	1:B:322:SER:HB2	1.81	0.63
1:B:82:ARG:CG	1:B:110:ARG:HB3	2.27	0.62
1:B:403:GLN:HE22	1:B:425:TRP:HE1	1.47	0.62
1:B:186:SER:HB3	1:B:189:ASP:HB2	1.82	0.62
1:A:366:MET:HB2	3:A:2176:HOH:O	2.00	0.61
1:B:280:GLU:HG3	1:B:311:PHE:CD2	2.36	0.60
1:A:69:ARG:HB3	1:A:70:PRO:CD	2.31	0.60
1:B:287:ALA:O	1:B:415:ALA:HA	2.03	0.59
1:A:219:ALA:O	1:A:223:ILE:HG13	2.04	0.58
1:A:135:LYS:HG3	3:A:2062:HOH:O	2.05	0.57
1:B:383:HIS:HE1	1:B:395:ARG:H	1.52	0.57
1:B:153:ALA:CB	1:B:176:HIS:HB2	2.31	0.57
1:A:132:ASP:HB3	1:A:135:LYS:HD3	1.86	0.57
1:A:318:HIS:HB3	1:B:306:ILE:HD13	1.86	0.56
1:A:306:ILE:HD12	1:B:308:GLN:HG2	1.87	0.56
1:B:181:LYS:HE2	1:B:265:ASP:O	2.06	0.56
1:B:83:ARG:HD3	1:B:389:ILE:CD1	2.36	0.56
1:B:352:LEU:HD12	1:B:364:SER:O	2.05	0.55
1:B:176:HIS:CD2	1:B:203:LYS:HB2	2.35	0.55
1:A:147:LYS:O	1:A:149:PRO:HD3	2.06	0.55
1:B:276:TYR:HA	1:B:433:TYR:CD2	2.42	0.54
1:B:174:GLY:O	1:B:175:LYS:HD3	2.07	0.54
1:B:375:ASN:ND2	1:B:377:PHE:HB2	2.22	0.54
1:B:100:LEU:N	1:B:101:PRO:HD2	2.22	0.54
1:B:87:ALA:HB3	1:B:154:VAL:HA	1.89	0.54
1:B:77:PRO:HD3	1:B:371:MET:O	2.07	0.54
1:B:306:ILE:HG12	1:B:320:ALA:HB2	1.89	0.54
1:A:287:ALA:O	1:A:415:ALA:HA	2.08	0.54
1:B:220:VAL:O	1:B:224:ARG:HG2	2.07	0.54
1:B:409:GLN:HG2	3:B:2080:HOH:O	2.09	0.53
1:B:83:ARG:HD3	1:B:389:ILE:HD13	1.89	0.53
1:B:428:VAL:HG13	1:B:430:GLN:HE22	1.72	0.53
1:B:374:ASN:HD22	1:B:375:ASN:H	1.57	0.53
1:B:171:PHE:CZ	1:B:177:VAL:HB	2.45	0.52
1:B:171:PHE:CZ	1:B:197:ALA:HB2	2.44	0.52
1:A:221:LYS:O	1:A:225:GLU:HG3	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:SER:HB3	1:B:122:ALA:HB2	1.91	0.52
1:B:96:LEU:HD23	1:B:100:LEU:HD12	1.92	0.52
1:B:276:TYR:CB	1:B:433:TYR:HB3	2.39	0.52
1:B:215:MET:CE	1:B:353:ILE:HG12	2.40	0.52
1:B:382:ASP:O	1:B:386:GLU:HB2	2.10	0.52
1:B:160:ASN:HD22	1:B:160:ASN:N	2.08	0.51
1:B:200:ALA:O	1:B:202:LYS:HG2	2.10	0.51
1:A:383:HIS:HE1	1:A:395:ARG:H	1.57	0.51
1:B:160:ASN:N	1:B:160:ASN:ND2	2.58	0.51
1:B:185:THR:HG23	1:B:259:GLY:HA3	1.93	0.51
1:A:290:TYR:HB3	1:B:316:LEU:HD11	1.93	0.50
1:B:375:ASN:HD21	1:B:377:PHE:HB2	1.76	0.50
1:B:211:HIS:HE1	1:B:272:ASN:ND2	1.99	0.50
1:B:187:VAL:O	1:B:190:CYS:HB2	2.11	0.50
1:A:238:ASN:HA	1:A:328:THR:O	2.10	0.50
1:B:215:MET:HE1	1:B:364:SER:HB2	1.94	0.50
1:A:383:HIS:CE1	1:A:395:ARG:H	2.28	0.50
1:B:63:PRO:C	1:B:65:THR:N	2.66	0.49
1:B:209:ARG:NH1	1:B:376:GLN:HE22	2.05	0.49
1:B:252:ARG:HA	1:B:258:ALA:HB2	1.94	0.49
1:B:403:GLN:NE2	1:B:403:GLN:HA	2.22	0.49
1:A:180:GLU:OE1	2:A:500:NDP:H2N	2.12	0.49
1:B:296:GLU:O	1:B:299:VAL:HG23	2.12	0.49
1:A:178:MET:SD	1:A:205:MET:HE3	2.52	0.49
1:A:261:GLY:N	1:A:302:GLU:OE1	2.45	0.49
1:B:222:LEU:HD12	1:B:222:LEU:C	2.32	0.49
1:B:286:ARG:HG3	1:B:286:ARG:HH11	1.78	0.48
1:B:238:ASN:HA	1:B:328:THR:O	2.12	0.48
1:B:91:LEU:HG	1:B:116:SER:HB2	1.95	0.48
1:A:181:LYS:HA	1:A:182:PRO:C	2.33	0.48
1:B:185:THR:HG22	1:B:185:THR:O	2.13	0.48
1:B:138:ASP:OD2	1:B:140:SER:HB2	2.14	0.48
1:A:124:ILE:O	1:A:128:GLU:HG3	2.14	0.48
1:B:404:ASP:O	1:B:408:ILE:HG13	2.15	0.47
1:B:413:GLU:HA	1:B:416:ARG:NH1	2.29	0.47
1:B:137:TYR:HE2	1:B:144:LYS:HB2	1.79	0.47
1:B:253:LEU:HD22	1:B:297:ARG:O	2.15	0.47
1:A:179:CYS:O	1:A:206:ILE:HA	2.14	0.47
1:B:276:TYR:HB2	1:B:433:TYR:HB3	1.96	0.47
1:A:89:VAL:HA	1:A:115:VAL:HB	1.95	0.47
1:A:76:ARG:HH11	1:A:76:ARG:HG2	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:ARG:HB3	1:A:70:PRO:HD2	1.97	0.46
1:A:211:HIS:HB2	1:A:212:TYR:CE2	2.50	0.46
1:B:173:ALA:O	1:B:175:LYS:HE2	2.15	0.46
1:B:298:PHE:CD2	1:B:303:ASP:HB2	2.50	0.46
1:B:176:HIS:CD2	1:B:203:LYS:H	2.33	0.46
1:A:286:ARG:NH2	1:B:288:TYR:CE1	2.84	0.45
1:A:116:SER:O	1:A:138:ASP:HA	2.16	0.45
1:A:219:ALA:HB1	1:A:340:LEU:HD21	1.99	0.45
1:A:181:LYS:HE3	2:A:500:NDP:N7N	2.31	0.45
1:B:403:GLN:CA	1:B:403:GLN:HE21	2.18	0.45
1:B:76:ARG:HD3	1:B:77:PRO:O	2.16	0.45
1:B:374:ASN:ND2	1:B:375:ASN:H	2.13	0.45
1:A:253:LEU:HD22	1:A:297:ARG:O	2.17	0.45
1:B:233:MET:HA	1:B:316:LEU:O	2.17	0.45
1:B:290:TYR:CD1	1:B:290:TYR:C	2.90	0.44
1:B:384:LEU:O	1:B:388:VAL:HG23	2.16	0.44
1:B:76:ARG:HG2	1:B:77:PRO:N	2.32	0.44
1:A:279:GLY:HA2	1:A:433:TYR:CD2	2.53	0.44
1:B:208:TYR:HB3	1:B:211:HIS:ND1	2.33	0.44
1:A:308:GLN:NE2	1:B:288:TYR:HB3	2.33	0.44
1:B:104:ALA:HB2	3:B:2013:HOH:O	2.18	0.44
1:B:417:THR:HB	1:B:419:ARG:HD3	1.99	0.44
1:B:298:PHE:CG	1:B:303:ASP:HB2	2.53	0.43
1:A:228:LEU:C	1:A:337:LYS:HB3	2.38	0.43
1:A:252:ARG:HH22	2:A:500:NDP:H6N	1.83	0.43
1:B:146:ALA:HA	1:B:175:LYS:HE2	2.00	0.43
1:B:235:THR:HB	1:B:332:SER:HB3	2.00	0.43
1:B:272:ASN:O	1:B:275:ARG:HB2	2.18	0.43
1:B:413:GLU:HG3	1:B:416:ARG:NH1	2.23	0.43
1:B:344:PRO:HB2	1:B:351:ASN:HA	2.00	0.43
1:B:87:ALA:HB3	1:B:154:VAL:HG23	2.01	0.43
1:A:273:GLY:HA2	1:A:276:TYR:CZ	2.53	0.43
1:B:181:LYS:HD2	1:B:181:LYS:C	2.39	0.43
1:B:69:ARG:HB3	1:B:70:PRO:HD2	2.01	0.43
1:A:182:PRO:HG3	1:A:264:MET:O	2.19	0.43
1:B:187:VAL:HG13	1:B:406:ARG:HG2	1.99	0.43
1:B:413:GLU:HA	1:B:416:ARG:CZ	2.49	0.43
1:A:332:SER:HA	1:A:340:LEU:O	2.18	0.42
1:B:189:ASP:O	1:B:193:MET:HG3	2.18	0.42
1:B:389:ILE:C	1:B:391:ASN:H	2.22	0.42
1:B:414:ALA:HB2	1:B:421:VAL:HG23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:307:TRP:CZ3	1:B:320:ALA:HA	2.54	0.42
1:B:178:MET:SD	1:B:205:MET:HE2	2.60	0.42
1:B:161:SER:HB3	1:B:257:LEU:HA	2.00	0.42
1:B:158:LEU:O	1:B:163:HIS:NE2	2.52	0.42
1:B:146:ALA:HB2	1:B:173:ALA:HA	2.01	0.42
1:B:235:THR:HA	1:B:318:HIS:O	2.19	0.42
1:A:334:GLN:NE2	1:B:325:THR:OG1	2.53	0.42
1:B:342:MET:O	1:B:345:ALA:HB2	2.20	0.42
1:A:267:GLY:CA	1:A:321:SER:HB3	2.50	0.42
1:B:95:ALA:O	1:B:100:LEU:HG	2.19	0.41
1:B:180:GLU:OE1	2:B:500:NDP:H2N	2.20	0.41
1:B:234:VAL:HA	1:B:332:SER:O	2.20	0.41
1:A:291:SER:O	1:A:292:ASP:C	2.59	0.41
1:B:100:LEU:HB3	1:B:129:TYR:CZ	2.54	0.41
1:B:160:ASN:H	1:B:160:ASN:ND2	2.17	0.41
1:A:326:THR:HG21	1:B:338:ALA:HA	2.02	0.41
1:A:368:GLN:N	1:A:368:GLN:NE2	2.46	0.41
1:A:94:TYR:CD1	2:A:500:NDP:H41N	2.56	0.41
1:B:375:ASN:ND2	1:B:378:SER:H	2.19	0.41
1:B:76:ARG:HH11	1:B:76:ARG:CG	2.34	0.41
1:B:164:ALA:HA	1:B:193:MET:SD	2.61	0.41
1:B:216:ASN:O	1:B:219:ALA:HB3	2.20	0.41
1:B:242:MET:SD	1:B:249:GLN:HB3	2.61	0.41
1:B:63:PRO:O	1:B:65:THR:N	2.54	0.41
1:A:214:PRO:HG3	1:A:373:ALA:HB3	2.03	0.41
1:A:226:ASN:HA	3:A:2108:HOH:O	2.20	0.40
1:B:183:MET:HE3	1:B:206:ILE:HD11	2.01	0.40
1:B:286:ARG:NH1	1:B:286:ARG:HG3	2.35	0.40
1:B:331:PHE:CD2	1:B:345:ALA:HB1	2.55	0.40
1:B:137:TYR:CE2	1:B:144:LYS:HB2	2.55	0.40
1:B:241:VAL:O	1:B:241:VAL:HG23	2.22	0.40
1:B:91:LEU:HB2	1:B:121:LYS:HE3	2.02	0.40
1:B:170:ALA:O	1:B:173:ALA:HB3	2.21	0.40
1:B:157:ILE:HG22	2:B:500:NDP:H51N	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	379/433 (88%)	360 (95%)	19 (5%)	0	100	100
1	B	379/433 (88%)	345 (91%)	30 (8%)	4 (1%)	14	26
All	All	758/866 (88%)	705 (93%)	49 (6%)	4 (0%)	29	48

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	64	THR
1	B	165	GLU
1	B	389	ILE
1	B	90	GLY

### 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	308/345 (89%)	299 (97%)	9 (3%)	42	69
1	B	308/345 (89%)	297 (96%)	11 (4%)	35	61
All	All	616/690 (89%)	596 (97%)	20 (3%)	39	65

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	ARG
1	A	78	MET

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Mol	Chain	Res	Type
1	A	118	ASN
1	A	155	TYR
1	A	181	LYS
1	A	240	ASP
1	A	244	GLN
1	A	368	GLN
1	A	374	ASN
1	B	76	ARG
1	B	181	LYS
1	B	204	LEU
1	B	240	ASP
1	B	244	GLN
1	B	308	GLN
1	B	349	TYR
1	B	374	ASN
1	B	384	LEU
1	B	403	GLN
1	B	430	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	GLN
1	A	97	ASN
1	A	118	ASN
1	A	226	ASN
1	A	227	GLN
1	A	244	GLN
1	A	250	GLN
1	A	308	GLN
1	A	334	GLN
1	A	368	GLN
1	A	374	ASN
1	A	383	HIS
1	A	390	ASN
1	A	409	GLN
1	B	98	GLN
1	B	160	ASN
1	B	176	HIS
1	B	211	HIS
1	B	226	ASN
1	B	227	GLN

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Mol	Chain	Res	Type
1	B	238	ASN
1	B	249	GLN
1	B	250	GLN
1	B	308	GLN
1	B	363	GLN
1	B	374	ASN
1	B	375	ASN
1	B	376	GLN
1	B	380	GLN
1	B	383	HIS
1	B	403	GLN
1	B	422	ASN
1	B	430	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

2 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	B	500	-	45,52,52	1.78	12 (26%)	53,80,80	1.27	5 (9%)



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NDP	A	500	-	45,52,52	1.78	11 (24%)	53,80,80	1.32	5 (9%)

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	B	500	-	-	9/30/77/77	0/5/5/5
2	NDP	A	500	-	-	7/30/77/77	0/5/5/5

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	NDP	C6N-C5N	4.93	1.42	1.33
2	B	500	NDP	P2B-O2X	-4.57	1.37	1.54
2	A	500	NDP	P2B-O2X	-4.49	1.37	1.54
2	B	500	NDP	C6N-C5N	4.45	1.41	1.33
2	B	500	NDP	P2B-O2B	4.13	1.67	1.59
2	A	500	NDP	O4B-C1B	3.56	1.46	1.41
2	A	500	NDP	C6N-N1N	3.28	1.45	1.37
2	B	500	NDP	C7N-C3N	3.09	1.55	1.48
2	A	500	NDP	P2B-O2B	3.03	1.65	1.59
2	B	500	NDP	C6N-N1N	2.99	1.44	1.37
2	B	500	NDP	C4N-C3N	-2.67	1.44	1.49
2	A	500	NDP	C4N-C3N	-2.67	1.44	1.49
2	A	500	NDP	C7N-C3N	2.63	1.54	1.48
2	B	500	NDP	O4B-C1B	2.57	1.44	1.41
2	B	500	NDP	C2N-C3N	2.48	1.41	1.34
2	A	500	NDP	C2N-C3N	2.47	1.41	1.34
2	B	500	NDP	C4A-N3A	2.32	1.38	1.35
2	B	500	NDP	P2B-O1X	2.22	1.57	1.50
2	A	500	NDP	C5A-N7A	-2.21	1.31	1.39
2	A	500	NDP	O4D-C1D	2.14	1.47	1.42
2	A	500	NDP	P2B-O1X	2.09	1.57	1.50
2	B	500	NDP	O4D-C1D	2.04	1.46	1.42
2	B	500	NDP	C2A-N3A	2.02	1.35	1.32

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	NDP	N3A-C2A-N1A	-5.14	120.64	128.68
2	B	500	NDP	N3A-C2A-N1A	-5.02	120.83	128.68
2	B	500	NDP	C3N-C2N-N1N	-3.52	118.07	123.10
2	A	500	NDP	C3N-C2N-N1N	-3.34	118.32	123.10
2	B	500	NDP	C1D-N1N-C2N	-2.54	116.88	121.11
2	A	500	NDP	C4A-C5A-N7A	2.48	111.98	109.40
2	A	500	NDP	C1D-N1N-C2N	-2.39	117.14	121.11
2	B	500	NDP	C4A-C5A-N7A	2.25	111.75	109.40
2	B	500	NDP	O2X-P2B-O2B	2.08	115.29	105.99
2	A	500	NDP	O2X-P2B-O2B	2.03	115.10	105.99

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	500	NDP	C5D-O5D-PN-O1N
2	B	500	NDP	C5D-O5D-PN-O2N
2	B	500	NDP	C5D-O5D-PN-O3
2	A	500	NDP	C5D-O5D-PN-O3
2	B	500	NDP	PA-O3-PN-O2N
2	A	500	NDP	PA-O3-PN-O2N
2	B	500	NDP	O4D-C1D-N1N-C2N
2	A	500	NDP	O4D-C1D-N1N-C2N
2	A	500	NDP	O4B-C4B-C5B-O5B
2	A	500	NDP	C2D-C1D-N1N-C2N
2	B	500	NDP	O4B-C4B-C5B-O5B
2	B	500	NDP	C2B-O2B-P2B-O3X
2	B	500	NDP	C2N-C3N-C7N-N7N
2	A	500	NDP	C5D-O5D-PN-O1N
2	A	500	NDP	C2N-C3N-C7N-N7N
2	B	500	NDP	C2D-C1D-N1N-C2N

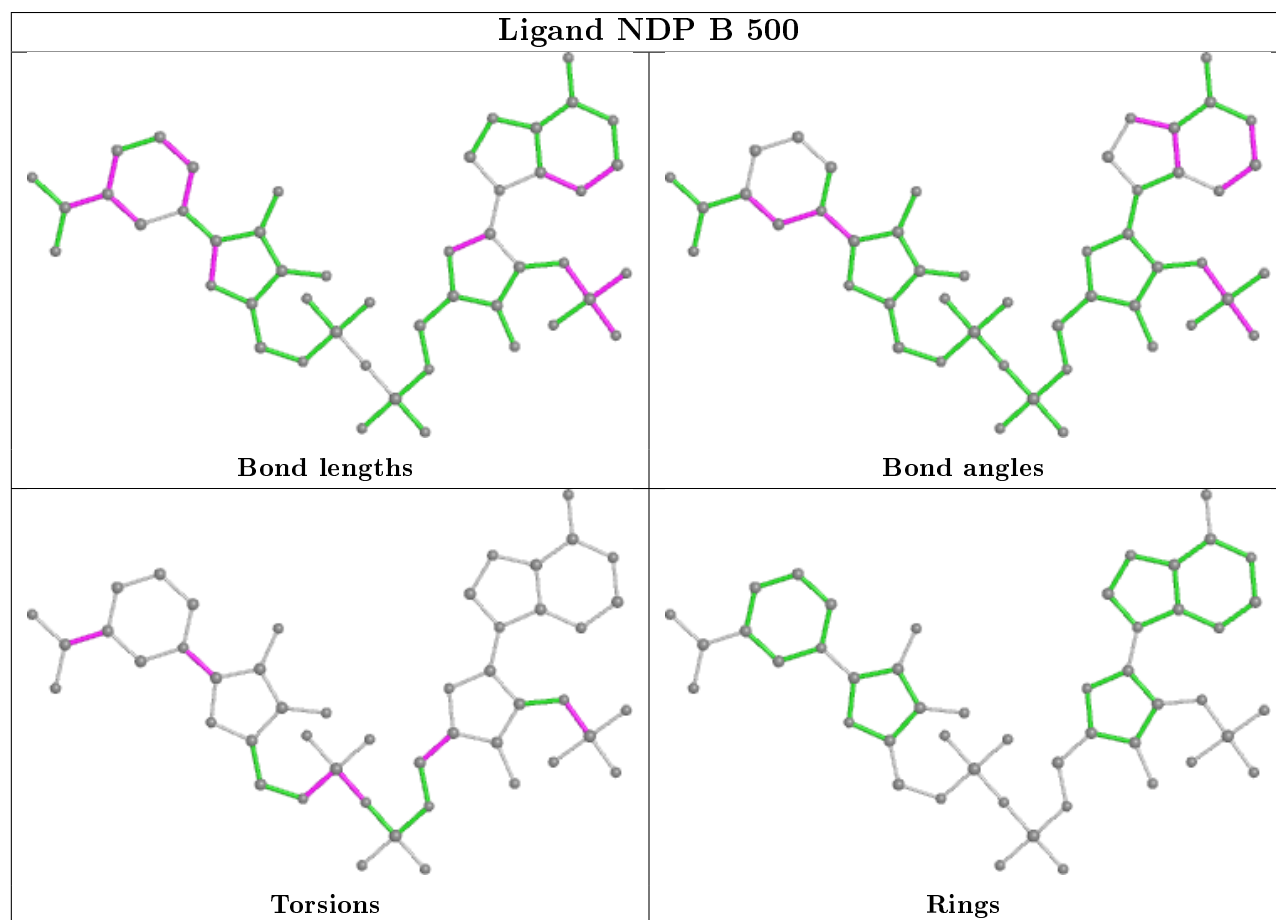
There are no ring outliers.

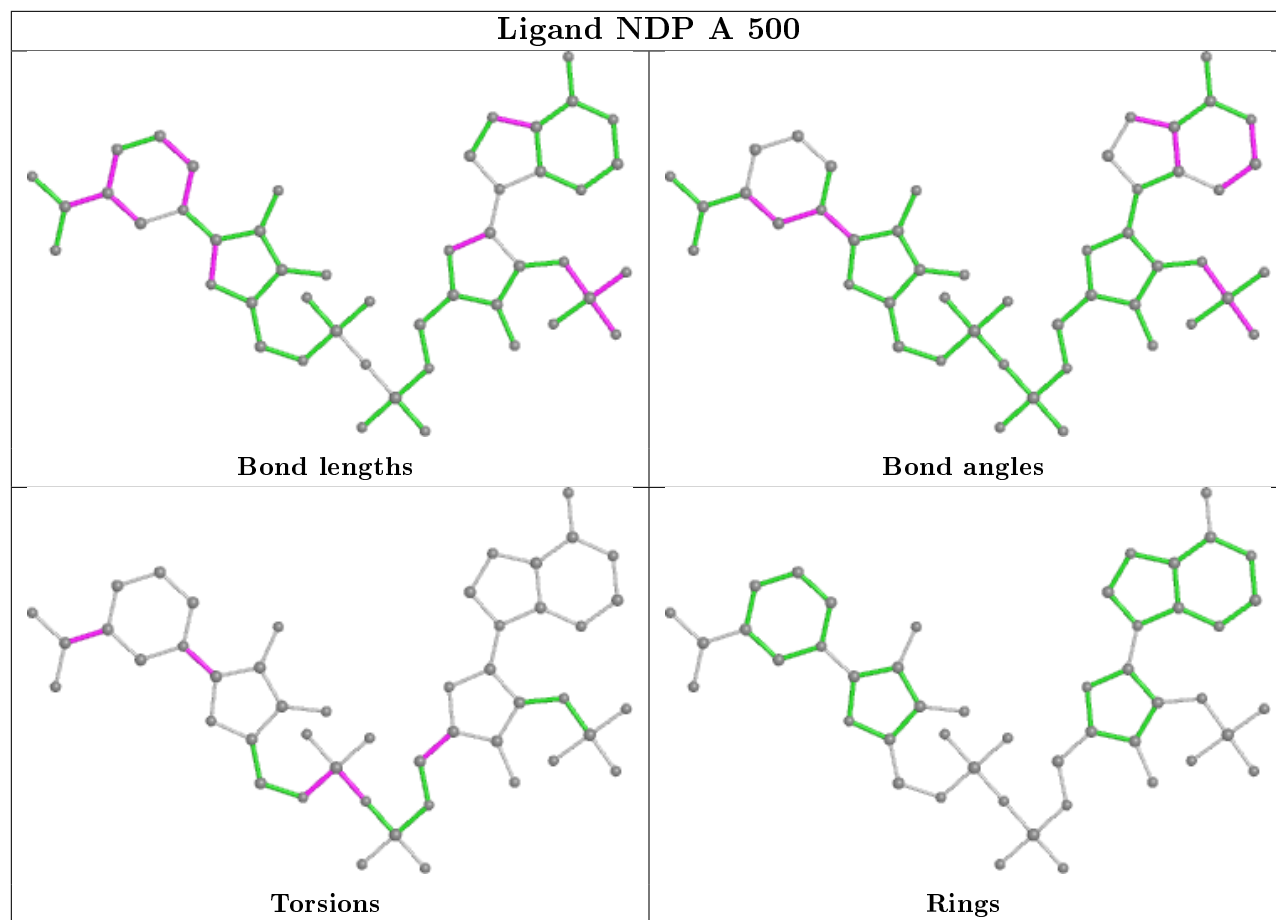
2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	500	NDP	2	0
2	A	500	NDP	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	381/433 (87%)	-0.28	2 (0%) 91 91	26, 36, 49, 63	0
1	B	381/433 (87%)	0.40	30 (7%) 12 12	31, 59, 82, 89	0
All	All	762/866 (87%)	0.06	32 (4%) 36 39	26, 43, 78, 89	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	266	ILE	4.1
1	B	430	GLN	3.5
1	B	428	VAL	3.4
1	B	269	TYR	3.4
1	B	419	ARG	3.2
1	B	267	GLY	3.2
1	B	181	LYS	3.1
1	B	265	ASP	3.0
1	B	236	THR	2.9
1	B	59	ALA	2.9
1	B	119	ALA	2.8
1	B	320	ALA	2.8
1	B	122	ALA	2.6
1	B	147	LYS	2.6
1	B	54	THR	2.6
1	B	134	ARG	2.5
1	B	141	ASN	2.5
1	B	53	ALA	2.4
1	B	239	SER	2.4
1	B	172	LYS	2.3
1	B	144	LYS	2.3
1	B	145	ILE	2.2
1	B	238	ASN	2.2
1	B	157	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	268	ILE	2.2
1	B	142	PHE	2.2
1	B	123	LYS	2.2
1	B	55	LEU	2.2
1	A	234	VAL	2.1
1	B	431	GLY	2.1
1	A	235	THR	2.1
1	B	182	PRO	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](#)

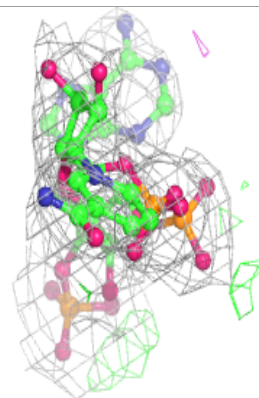
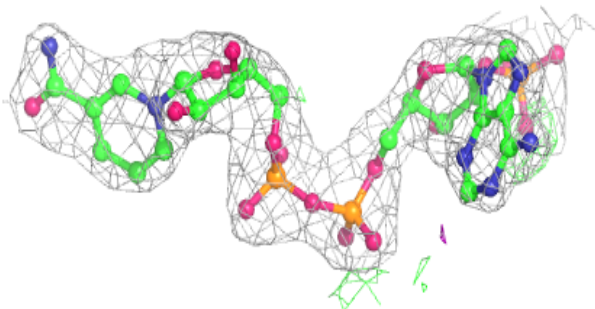
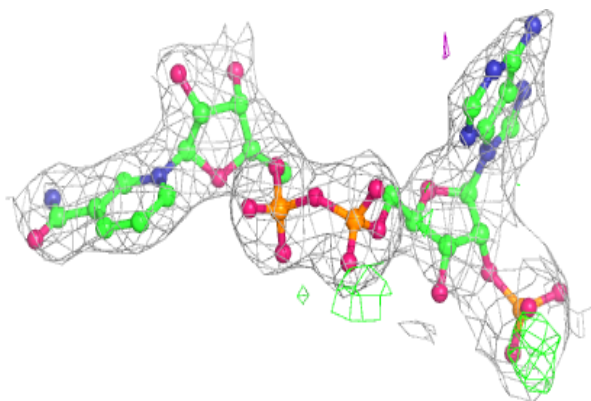
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	NDP	B	500	48/48	0.89	0.21	59,64,69,70	0
2	NDP	A	500	48/48	0.97	0.18	29,35,41,43	0

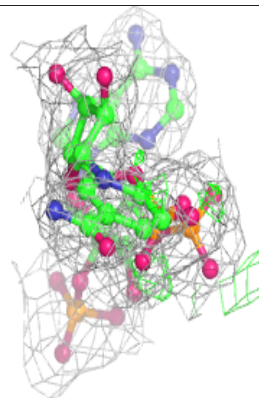
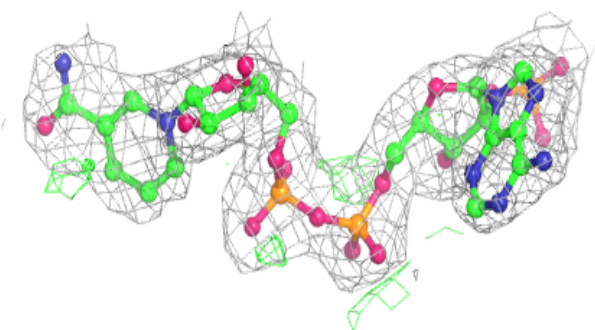
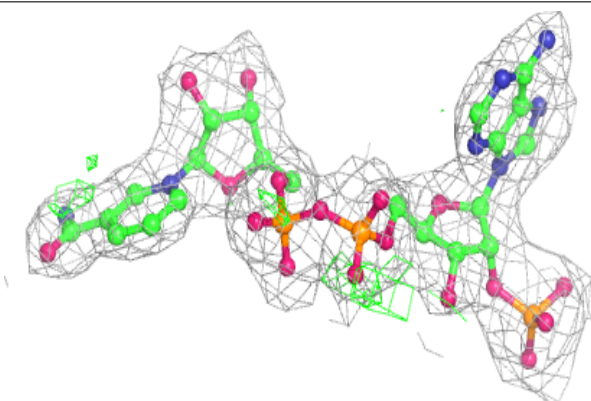
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around NDP B 500:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around NDP A 500:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.