



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 09:40 am GMT

PDB ID : 1N10  
Title : Crystal Structure of Phl p 1, a Major Timothy Grass Pollen Allergen  
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Deposited on : 2002-10-16  
Resolution : 2.90 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

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

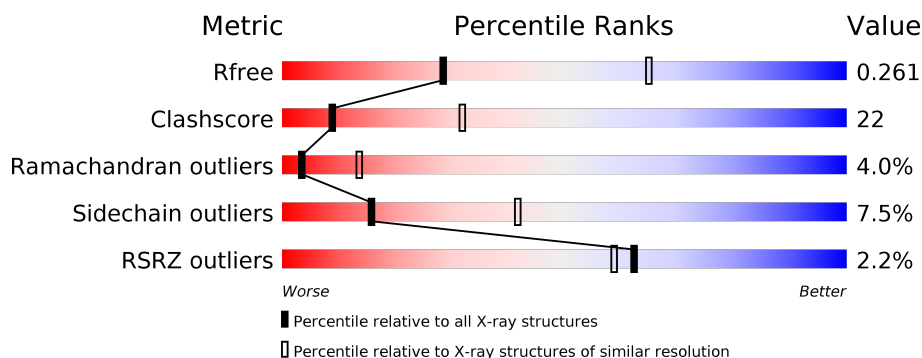
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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))
$R_{free}$	100719	1586 (2.90-2.90)
Clashscore	112137	1807 (2.90-2.90)
Ramachandran outliers	110173	1768 (2.90-2.90)
Sidechain outliers	110143	1770 (2.90-2.90)
RSRZ outliers	101464	1596 (2.90-2.90)

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. 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	241	<div> <div>2%</div> <div> <div></div> <div>55%</div> <div>34%</div> <div>5% • 5%</div> </div> </div>
1	B	241	<div> <div>2%</div> <div> <div></div> <div>55%</div> <div>34%</div> <div>5% • 5%</div> </div> </div>

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 3540 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.

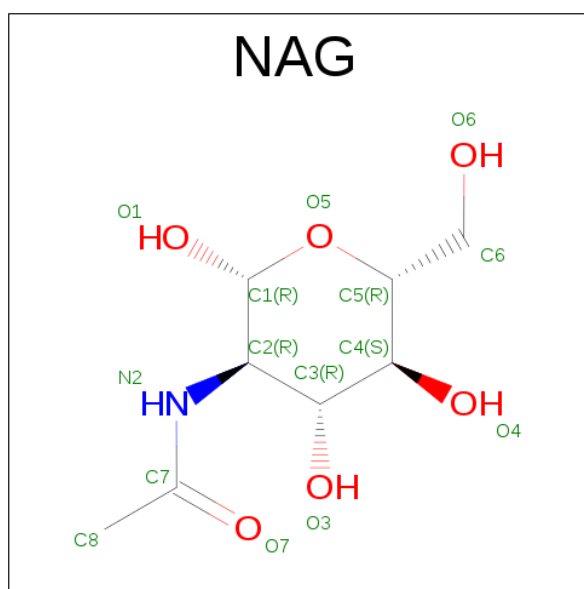
- Molecule 1 is a protein called Pollen allergen Phl p 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	0	0	0
			1756	1114	292	341	9			
1	B	228	Total	C	N	O	S	0	0	0
			1756	1114	292	341	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1001	ALA	GLY	SEE REMARK 999	UNP P43213
B	2001	ALA	GLY	SEE REMARK 999	UNP P43213

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



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

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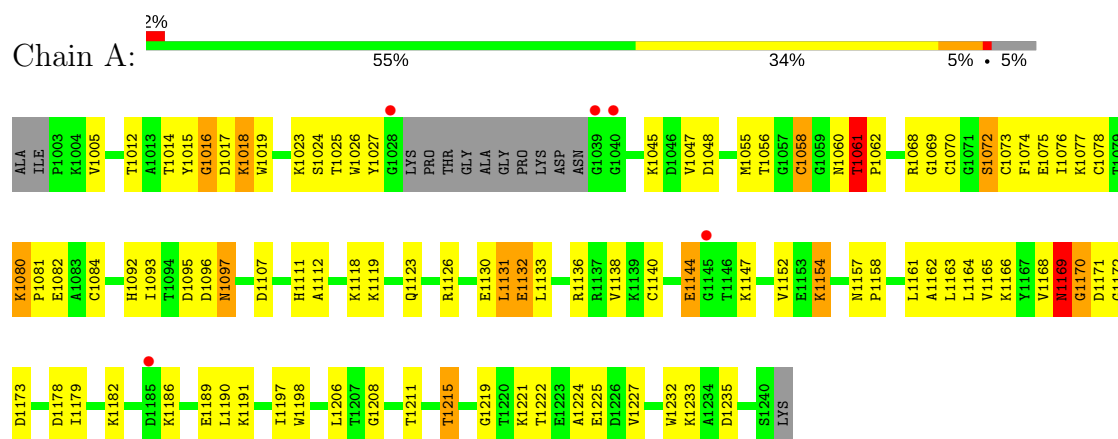
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

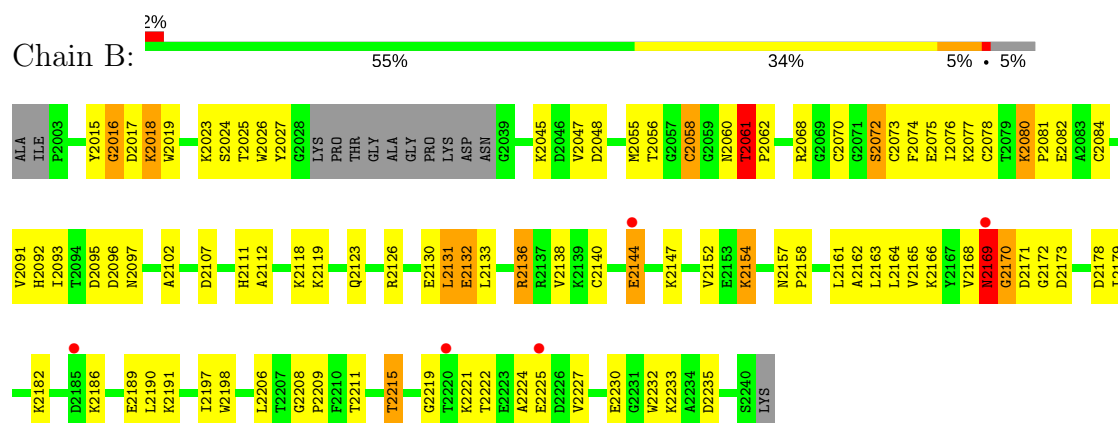
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Pollen allergen Phl p 1



- Molecule 1: Pollen allergen Phl p 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	59.69Å 65.53Å 144.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.90 19.90 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.6 (20.00-2.90) 98.6 (19.90-2.90)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	21.24 (at 2.88Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.239 , 0.267 0.240 , 0.261	Depositor DCC
$R_{free}$ test set	664 reflections (5.13%)	DCC
Wilson B-factor (Å <sup>2</sup> )	50.2	Xtriage
Anisotropy	0.515	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 31.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	3540	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.00% 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: NAG

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	0.43	1/1803 (0.1%)	0.71	1/2440 (0.0%)
1	B	0.41	0/1803	0.70	1/2440 (0.0%)
All	All	0.42	1/3606 (0.0%)	0.70	2/4880 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1058	CYS	CB-SG	6.10	1.92	1.82

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1058	CYS	CA-CB-SG	5.60	124.08	114.00
1	B	2058	CYS	CA-CB-SG	5.17	123.30	114.00

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	1756	0	1692	79	0
1	B	1756	0	1692	81	0
2	A	14	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	14	0	13	0	0
All	All	3540	0	3410	156	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 22.

All (156) 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:2027:TYR:HE2	1:B:2058:CYS:HG	1.18	0.92
1:B:2027:TYR:CE2	1:B:2058:CYS:SG	2.66	0.89
1:A:1027:TYR:CE2	1:A:1058:CYS:SG	2.76	0.78
1:B:2075:GLU:HG3	1:B:2138:VAL:HG11	1.65	0.78
1:A:1075:GLU:HG3	1:A:1138:VAL:HG11	1.66	0.78
1:A:1027:TYR:HE2	1:A:1058:CYS:HG	1.32	0.76
1:A:1224:ALA:HB1	1:A:1227:VAL:HG21	1.68	0.76
1:B:2224:ALA:HB1	1:B:2227:VAL:HG21	1.67	0.76
1:A:1058:CYS:HB2	1:A:1107:ASP:HB3	1.69	0.74
1:B:2027:TYR:HE2	1:B:2058:CYS:SG	2.05	0.73
1:A:1026:TRP:CE2	1:A:1126:ARG:HG2	2.25	0.71
1:B:2058:CYS:HB2	1:B:2107:ASP:HB3	1.72	0.70
1:B:2026:TRP:CE2	1:B:2126:ARG:HG2	2.27	0.70
1:A:1026:TRP:CZ2	1:A:1126:ARG:HG2	2.26	0.70
1:A:1157:ASN:HB2	1:A:1158:PRO:HD2	1.75	0.68
1:B:2157:ASN:HB2	1:B:2158:PRO:HD2	1.74	0.68
1:A:1025:THR:HG22	1:A:1130:GLU:HG2	1.76	0.68
1:A:1027:TYR:HE2	1:A:1058:CYS:SG	2.17	0.67
1:A:1080:LYS:HB3	1:A:1081:PRO:CD	2.24	0.67
1:B:2026:TRP:CZ2	1:B:2126:ARG:HG2	2.30	0.66
1:B:2080:LYS:HB3	1:B:2081:PRO:CD	2.25	0.66
1:B:2024:SER:HB3	1:B:2131:LEU:HD23	1.78	0.65
1:B:2025:THR:HG22	1:B:2130:GLU:HG2	1.79	0.63
1:A:1024:SER:HB3	1:A:1131:LEU:HD23	1.80	0.62
1:B:2164:LEU:HD21	1:B:2197:ILE:HG12	1.81	0.61
1:A:1123:GLN:HA	1:A:1126:ARG:NH2	2.16	0.60
1:B:2123:GLN:HA	1:B:2126:ARG:NH2	2.17	0.60
1:B:2024:SER:HB3	1:B:2131:LEU:CD2	2.32	0.59
1:B:2190:LEU:HB3	1:B:2198:TRP:HB3	1.84	0.59
1:A:1215:THR:CG2	1:A:1221:LYS:HG2	2.32	0.59
1:A:1164:LEU:HD21	1:A:1197:ILE:HG12	1.83	0.59
1:A:1024:SER:HB3	1:A:1131:LEU:CD2	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1012:THR:O	1:A:1014:THR:HG23	2.04	0.57
1:A:1061:THR:HG23	1:A:1062:PRO:HD3	1.85	0.57
1:A:1154:LYS:HD2	1:A:1154:LYS:H	1.69	0.57
1:B:2055:MET:HE2	1:B:2112:ALA:HA	1.87	0.56
1:A:1190:LEU:HB3	1:A:1198:TRP:HB3	1.86	0.56
1:A:1055:MET:HE2	1:A:1112:ALA:HA	1.87	0.56
1:B:2215:THR:CG2	1:B:2221:LYS:HG2	2.35	0.56
1:A:1023:LYS:HD2	1:A:1132:GLU:OE1	2.05	0.56
1:A:1224:ALA:CB	1:A:1227:VAL:HG21	2.37	0.55
1:B:2224:ALA:HB1	1:B:2227:VAL:CG2	2.37	0.54
1:B:2154:LYS:HD2	1:B:2154:LYS:H	1.72	0.54
1:A:1224:ALA:HB1	1:A:1227:VAL:CG2	2.37	0.54
1:B:2080:LYS:CB	1:B:2081:PRO:CD	2.86	0.54
1:B:2023:LYS:HD2	1:B:2132:GLU:OE1	2.08	0.54
1:A:1080:LYS:CB	1:A:1081:PRO:CD	2.86	0.53
1:B:2061:THR:HG23	1:B:2062:PRO:HD3	1.90	0.53
1:A:1144:GLU:OE1	1:A:1144:GLU:N	2.42	0.53
1:A:1163:LEU:HD21	1:A:1190:LEU:CD2	2.39	0.53
1:B:2232:TRP:C	1:B:2233:LYS:HD3	2.28	0.53
1:A:1232:TRP:C	1:A:1233:LYS:HD3	2.29	0.52
1:A:1047:VAL:HG13	1:A:1056:THR:HG21	1.92	0.52
1:A:1224:ALA:O	1:A:1227:VAL:HG23	2.10	0.52
1:B:2047:VAL:HG13	1:B:2056:THR:HG21	1.91	0.52
1:B:2080:LYS:HD3	1:B:2131:LEU:HB2	1.92	0.52
1:B:2224:ALA:O	1:B:2227:VAL:HG23	2.10	0.52
1:B:2015:TYR:O	1:B:2016:GLY:O	2.28	0.51
1:B:2045:LYS:HZ3	1:B:2166:LYS:HZ1	1.57	0.51
1:B:2168:VAL:C	1:B:2170:GLY:H	2.13	0.51
1:B:2018:LYS:HB3	1:B:2018:LYS:NZ	2.25	0.51
1:B:2025:THR:CG2	1:B:2130:GLU:HG2	2.41	0.51
1:A:1018:LYS:HB3	1:A:1018:LYS:NZ	2.26	0.51
1:B:2179:ILE:HA	1:B:2211:THR:O	2.11	0.51
1:A:1025:THR:CG2	1:A:1130:GLU:HG2	2.41	0.50
1:B:2163:LEU:HD21	1:B:2190:LEU:CD2	2.41	0.50
1:A:1168:VAL:C	1:A:1170:GLY:H	2.14	0.50
1:A:1078:CYS:O	1:A:1084:CYS:HB3	2.11	0.50
1:B:2224:ALA:CB	1:B:2227:VAL:HG21	2.38	0.49
1:B:2144:GLU:N	1:B:2144:GLU:OE1	2.46	0.49
1:B:2078:CYS:O	1:B:2084:CYS:HB3	2.13	0.49
1:B:2080:LYS:CD	1:B:2132:GLU:H	2.26	0.48
1:A:1186:LYS:NZ	1:B:2186:LYS:HB3	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2152:VAL:O	1:B:2235:ASP:N	2.46	0.48
1:B:2191:LYS:HD2	1:B:2191:LYS:N	2.29	0.48
1:A:1080:LYS:HD3	1:A:1131:LEU:HB2	1.94	0.48
1:A:1179:ILE:HA	1:A:1211:THR:O	2.14	0.47
1:A:1015:TYR:O	1:A:1016:GLY:O	2.32	0.47
1:A:1161:LEU:HB2	1:A:1206:LEU:HD13	1.95	0.47
1:A:1161:LEU:CD1	1:A:1163:LEU:HD22	2.45	0.47
1:A:1163:LEU:HD23	1:A:1163:LEU:H	1.80	0.47
1:B:2068:ARG:HG3	1:B:2173:ASP:HB3	1.96	0.47
1:A:1152:VAL:O	1:A:1235:ASP:N	2.46	0.47
1:A:1140:CYS:HB2	1:A:1169:ASN:ND2	2.30	0.47
1:B:2080:LYS:HD2	1:B:2132:GLU:H	1.79	0.47
1:A:1005:VAL:HG21	1:B:2102:ALA:HA	1.97	0.47
1:A:1169:ASN:O	1:A:1170:GLY:O	2.32	0.47
1:B:2171:ASP:OD1	1:B:2172:GLY:N	2.48	0.46
1:A:1045:LYS:HZ3	1:A:1166:LYS:HZ1	1.62	0.46
1:B:2169:ASN:O	1:B:2170:GLY:O	2.33	0.46
1:A:1191:LYS:HD2	1:A:1191:LYS:N	2.30	0.46
1:A:1080:LYS:CD	1:A:1132:GLU:H	2.28	0.46
1:B:2161:LEU:CD1	1:B:2163:LEU:HD22	2.45	0.46
1:A:1080:LYS:HD2	1:A:1132:GLU:H	1.80	0.46
1:B:2075:GLU:HG3	1:B:2138:VAL:CG1	2.39	0.46
1:B:2161:LEU:HB2	1:B:2206:LEU:HD13	1.96	0.46
1:A:1215:THR:HG22	1:A:1221:LYS:HG2	1.95	0.46
1:B:2055:MET:HA	1:B:2055:MET:CE	2.45	0.46
1:B:2072:SER:HB3	1:B:2074:PHE:CE1	2.51	0.46
1:A:1133:LEU:C	1:A:1133:LEU:HD12	2.36	0.46
1:A:1080:LYS:HB3	1:A:1081:PRO:HD3	1.98	0.46
1:B:2123:GLN:HA	1:B:2126:ARG:HH22	1.81	0.46
1:B:2140:CYS:HB2	1:B:2169:ASN:ND2	2.30	0.46
1:B:2073:CYS:SG	1:B:2092:HIS:CD2	3.08	0.45
1:B:2080:LYS:HB3	1:B:2081:PRO:HD2	1.98	0.45
1:A:1123:GLN:HA	1:A:1126:ARG:HH22	1.79	0.45
1:B:2163:LEU:H	1:B:2163:LEU:HD23	1.81	0.45
1:A:1075:GLU:HG3	1:A:1138:VAL:CG1	2.41	0.45
1:B:2211:THR:OG1	1:B:2225:GLU:HA	2.16	0.45
1:A:1072:SER:HB3	1:A:1074:PHE:CE1	2.51	0.45
1:A:1186:LYS:HD2	1:B:2186:LYS:HD2	1.99	0.45
1:A:1171:ASP:OD1	1:A:1172:GLY:N	2.49	0.45
1:A:1055:MET:HA	1:A:1055:MET:CE	2.47	0.45
1:A:1165:VAL:HG12	1:A:1168:VAL:CG2	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1178:ASP:OD1	1:A:1189:GLU:HG2	2.17	0.44
1:A:1068:ARG:HG3	1:A:1173:ASP:HB3	2.00	0.44
1:A:1005:VAL:HG21	1:B:2102:ALA:CA	2.48	0.44
1:A:1070:CYS:O	1:A:1093:ILE:O	2.36	0.44
1:A:1211:THR:OG1	1:A:1225:GLU:HA	2.17	0.44
1:B:2027:TYR:OH	1:B:2058:CYS:SG	2.74	0.43
1:A:1073:CYS:SG	1:A:1092:HIS:CD2	3.12	0.43
1:A:1019:TRP:CZ2	1:A:1077:LYS:HE3	2.54	0.43
1:A:1095:ASP:CG	1:A:1096:ASP:H	2.22	0.43
1:A:1047:VAL:CG1	1:A:1056:THR:HG21	2.47	0.43
1:B:2060:ASN:ND2	1:B:2062:PRO:HD2	2.34	0.43
1:B:2133:LEU:C	1:B:2133:LEU:HD12	2.39	0.43
1:B:2082:GLU:OE2	1:B:2082:GLU:HA	2.19	0.43
1:B:2019:TRP:CZ2	1:B:2077:LYS:HE3	2.55	0.42
1:B:2080:LYS:HB3	1:B:2081:PRO:HD3	2.01	0.42
1:B:2076:ILE:O	1:B:2076:ILE:HG13	2.19	0.42
1:B:2047:VAL:CG1	1:B:2056:THR:HG21	2.48	0.42
1:B:2164:LEU:CD2	1:B:2197:ILE:HG12	2.49	0.42
1:B:2215:THR:HG22	1:B:2221:LYS:HG2	2.00	0.42
1:A:1055:MET:HE3	1:A:1111:HIS:ND1	2.34	0.42
1:B:2165:VAL:HG12	1:B:2168:VAL:CG2	2.50	0.42
1:A:1082:GLU:HA	1:A:1082:GLU:OE2	2.19	0.42
1:B:2070:CYS:O	1:B:2093:ILE:O	2.38	0.42
1:B:2095:ASP:CG	1:B:2096:ASP:H	2.23	0.42
1:B:2060:ASN:CG	1:B:2062:PRO:HD2	2.39	0.41
1:A:1118:LYS:O	1:A:1119:LYS:C	2.59	0.41
1:A:1162:ALA:HA	1:A:1198:TRP:O	2.20	0.41
1:A:1164:LEU:CD2	1:A:1197:ILE:HG12	2.51	0.41
1:B:2073:CYS:HA	1:B:2091:VAL:O	2.20	0.41
1:A:1061:THR:HG23	1:A:1062:PRO:CD	2.47	0.41
1:A:1069:GLY:O	1:A:1072:SER:HB2	2.21	0.41
1:A:1163:LEU:HD21	1:A:1190:LEU:HD23	2.03	0.41
1:A:1097:ASN:C	1:A:1097:ASN:HD22	2.25	0.41
1:B:2178:ASP:OD1	1:B:2189:GLU:HG2	2.21	0.41
1:B:2162:ALA:HA	1:B:2198:TRP:O	2.21	0.41
1:B:2047:VAL:CG1	1:B:2047:VAL:O	2.69	0.40
1:A:1060:ASN:ND2	1:A:1062:PRO:HD2	2.36	0.40
1:B:2075:GLU:OE1	1:B:2136:ARG:NH2	2.52	0.40
1:B:2055:MET:HE3	1:B:2111:HIS:ND1	2.36	0.40
1:B:2118:LYS:O	1:B:2119:LYS:C	2.59	0.40
1:A:1076:ILE:HG13	1:A:1076:ILE:O	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2209:PRO:HD3	1:B:2230:GLU:HB2	2.03	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	224/241 (93%)	197 (88%)	18 (8%)	9 (4%)	3	14
1	B	224/241 (93%)	199 (89%)	16 (7%)	9 (4%)	3	14
All	All	448/482 (93%)	396 (88%)	34 (8%)	18 (4%)	3	14

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1080	LYS
1	B	2080	LYS
1	A	1016	GLY
1	A	1170	GLY
1	A	1182	LYS
1	B	2016	GLY
1	B	2170	GLY
1	B	2182	LYS
1	B	2072	SER
1	B	2169	ASN
1	A	1072	SER
1	A	1169	ASN
1	A	1208	GLY
1	A	1061	THR
1	A	1219	GLY
1	B	2061	THR
1	B	2208	GLY

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Mol	Chain	Res	Type
1	B	2219	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	187/196 (95%)	173 (92%)	14 (8%)	16	42
1	B	187/196 (95%)	173 (92%)	14 (8%)	16	42
All	All	374/392 (95%)	346 (92%)	28 (8%)	16	42

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1017	ASP
1	A	1018	LYS
1	A	1048	ASP
1	A	1061	THR
1	A	1097	ASN
1	A	1131	LEU
1	A	1132	GLU
1	A	1136	ARG
1	A	1144	GLU
1	A	1147	LYS
1	A	1154	LYS
1	A	1169	ASN
1	A	1215	THR
1	A	1222	THR
1	B	2017	ASP
1	B	2018	LYS
1	B	2048	ASP
1	B	2061	THR
1	B	2097	ASN
1	B	2131	LEU
1	B	2132	GLU
1	B	2136	ARG

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Mol	Chain	Res	Type
1	B	2144	GLU
1	B	2147	LYS
1	B	2154	LYS
1	B	2169	ASN
1	B	2215	THR
1	B	2222	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1092	HIS
1	A	1097	ASN
1	A	1169	ASN
1	B	2092	HIS
1	B	2097	ASN
1	B	2169	ASN

### 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	NAG	A	3001	1	14,14,15	1.28	4 (28%)	15,19,21	1.34	1 (6%)
2	NAG	B	3002	1	14,14,15	1.28	2 (14%)	15,19,21	1.34	1 (6%)

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	NAG	A	3001	1	-	0/6/23/26	0/1/1/1
2	NAG	B	3002	1	-	0/6/23/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	3001	NAG	O3-C3	2.09	1.47	1.43
2	A	3001	NAG	C4-C5	2.13	1.57	1.53
2	A	3001	NAG	O5-C5	2.15	1.48	1.43
2	A	3001	NAG	C2-N2	2.18	1.50	1.46
2	B	3002	NAG	C4-C5	2.30	1.57	1.53
2	B	3002	NAG	O5-C5	2.52	1.48	1.43

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	3002	NAG	C1-O5-C5	3.19	116.57	112.17
2	A	3001	NAG	C1-O5-C5	3.59	117.12	112.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	228/241 (94%)	-0.25	5 (2%) 62 59	18, 38, 76, 100	0
1	B	228/241 (94%)	-0.19	5 (2%) 62 59	19, 42, 85, 114	0
All	All	456/482 (94%)	-0.22	10 (2%) 62 59	18, 40, 82, 114	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1039	GLY	4.6
1	A	1145	GLY	3.0
1	B	2169	ASN	2.7
1	A	1040	GLY	2.6
1	B	2144	GLU	2.6
1	B	2220	THR	2.4
1	B	2225	GLU	2.2
1	A	1028	GLY	2.2
1	B	2185	ASP	2.2
1	A	1185	ASP	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, 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NAG	B	3002	14/15	0.79	0.29	-	47,53,63,68	0
2	NAG	A	3001	14/15	0.82	0.34	-	47,51,63,70	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.