



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:37 PM GMT

PDB ID : 1GIC
Title : CONCAVALIN A COMPLEXED WITH METHYL ALPHA-D-GLUCOPYRANOSIDE
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Deposited on : 1996-08-08
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (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

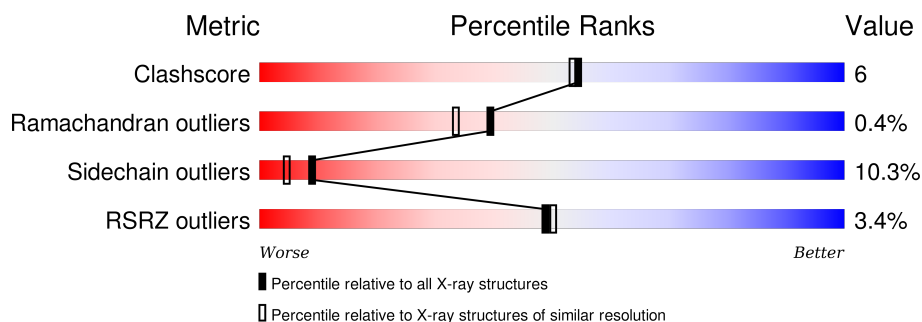
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	237	<div> <div>4%</div> <div>77%</div> <div>21%</div> <div>..</div> </div>
1	B	237	<div> <div>3%</div> <div>74%</div> <div>19%</div> <div>6%</div> <div>.</div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4008 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.

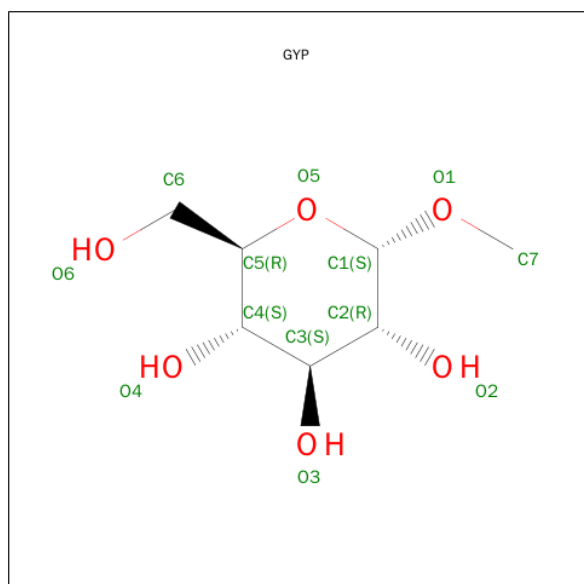
- Molecule 1 is a protein called CONCANAVALIN A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	0	0	0
			1809	1141	302	364	2			
1	B	237	Total	C	N	O	S	0	0	0
			1809	1141	302	364	2			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	151	ASP	GLU	CONFLICT	UNP P02866
A	155	GLU	ARG	CONFLICT	UNP P02866
B	151	ASP	GLU	CONFLICT	UNP P02866
B	155	GLU	ARG	CONFLICT	UNP P02866

- Molecule 2 is SUGAR (O1-METHYL-GLUCOSE) (three-letter code: GYP) (formula: C₇H₁₄O₆).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 13 7 6	0	0
2	B	1	Total C O 13 7 6	0	0

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Mn 1 1	0	0
3	A	1	Total Mn 1 1	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Ca 1 1	0	0
4	A	1	Total Ca 1 1	0	0

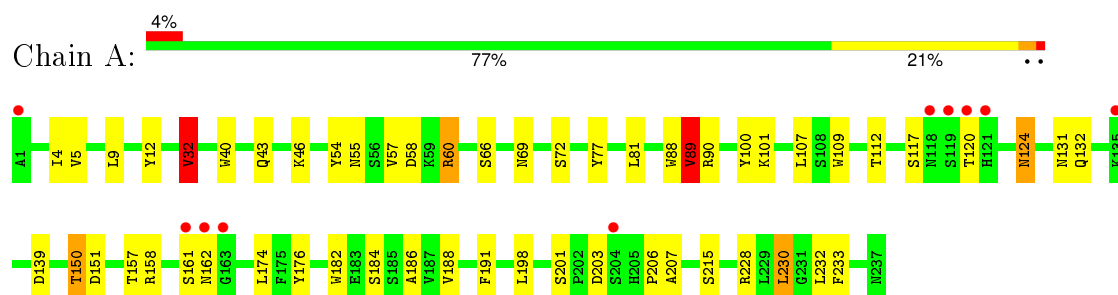
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	186	Total O 186 186	0	0
5	B	174	Total O 174 174	0	0

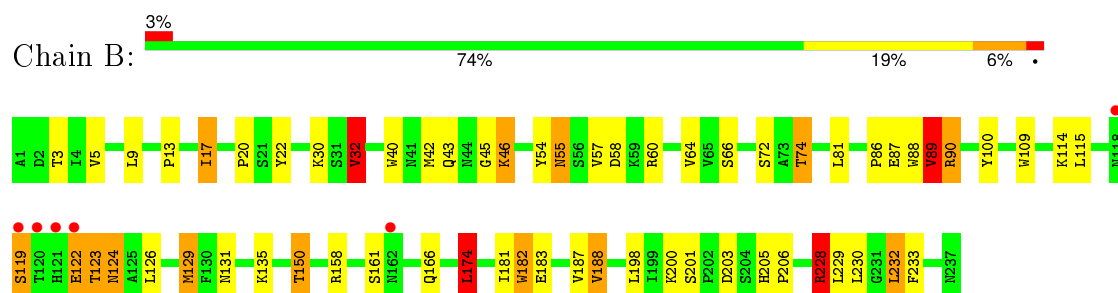
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: CONCANAVALIN A



• Molecule 1: CONCANAVALIN A



4 Data and refinement statistics

Property	Value	Source
Space group	I 21 3	Depositor
Cell constants a, b, c, α , β , γ	167.80Å 167.80Å 167.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.00 – 2.00 44.85 – 2.02	Depositor EDS
% Data completeness (in resolution range)	64.0 (100.00-2.00) 63.6 (44.85-2.02)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.68 (at 2.01Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.170 , (Not available) 0.176 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	19.8	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 70.1	EDS
Estimated twinning fraction	0.020 for -l,-k,-h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 34663 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4008	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.26% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 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: CA, MN, GYP

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	0.89	1/1851 (0.1%)	1.58	25/2522 (1.0%)
1	B	0.93	1/1851 (0.1%)	1.69	27/2522 (1.1%)
All	All	0.91	2/3702 (0.1%)	1.63	52/5044 (1.0%)

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	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	188	VAL	CB-CG2	-5.24	1.41	1.52
1	A	188	VAL	CB-CG2	-5.17	1.42	1.52

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	90	ARG	NE-CZ-NH2	-12.57	114.02	120.30
1	B	228	ARG	NE-CZ-NH1	12.10	126.35	120.30
1	B	90	ARG	NE-CZ-NH1	11.50	126.05	120.30
1	B	22	TYR	CB-CG-CD2	-9.91	115.05	121.00
1	A	89	VAL	N-CA-CB	-8.83	92.08	111.50
1	A	182	TRP	CD1-CG-CD2	8.74	113.30	106.30
1	A	228	ARG	NE-CZ-NH1	8.57	124.59	120.30
1	B	89	VAL	N-CA-CB	-8.38	93.07	111.50
1	A	228	ARG	NE-CZ-NH2	-8.34	116.13	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	88	TRP	CD1-CG-CD2	7.94	112.65	106.30
1	A	182	TRP	CE2-CD2-CG	-7.70	101.14	107.30
1	B	88	TRP	CD1-CG-CD2	7.56	112.35	106.30
1	B	40	TRP	CD1-CG-CD2	7.42	112.24	106.30
1	B	174	LEU	CA-CB-CG	7.35	132.21	115.30
1	A	88	TRP	CE2-CD2-CG	-7.28	101.47	107.30
1	B	32	VAL	CB-CA-C	-7.17	97.77	111.40
1	A	109	TRP	CD1-CG-CD2	7.15	112.02	106.30
1	A	40	TRP	CD1-CG-CD2	7.08	111.97	106.30
1	A	40	TRP	CE2-CD2-CG	-7.01	101.69	107.30
1	A	77	TYR	CB-CG-CD2	-6.96	116.83	121.00
1	B	40	TRP	CE2-CD2-CG	-6.91	101.78	107.30
1	A	32	VAL	CB-CA-C	-6.88	98.32	111.40
1	B	88	TRP	CE2-CD2-CG	-6.85	101.82	107.30
1	B	182	TRP	CD1-CG-CD2	6.82	111.76	106.30
1	B	228	ARG	NE-CZ-NH2	-6.78	116.91	120.30
1	B	182	TRP	CE2-CD2-CG	-6.77	101.88	107.30
1	A	12	TYR	CB-CG-CD2	-6.62	117.03	121.00
1	B	232	LEU	CA-CB-CG	6.18	129.52	115.30
1	B	109	TRP	CE2-CD2-CG	-6.13	102.40	107.30
1	A	182	TRP	CG-CD1-NE1	-6.05	104.05	110.10
1	A	182	TRP	CA-CB-CG	6.02	125.14	113.70
1	B	109	TRP	CD1-CG-CD2	5.99	111.09	106.30
1	B	42	MET	CG-SD-CE	-5.88	90.78	100.20
1	B	5	VAL	N-CA-C	-5.87	95.15	111.00
1	A	88	TRP	CG-CD2-CE3	5.85	139.16	133.90
1	A	109	TRP	CE2-CD2-CG	-5.82	102.65	107.30
1	B	22	TYR	CB-CG-CD1	5.81	124.48	121.00
1	A	89	VAL	CG1-CB-CG2	5.79	120.16	110.90
1	B	114	LYS	CB-CG-CD	-5.65	96.92	111.60
1	A	54	TYR	CB-CG-CD1	-5.58	117.65	121.00
1	A	88	TRP	CB-CG-CD1	-5.47	119.89	127.00
1	A	9	LEU	N-CA-C	-5.32	96.63	111.00
1	A	230	LEU	CA-CB-CG	5.32	127.54	115.30
1	B	17	ILE	CA-CB-CG1	-5.31	100.91	111.00
1	B	90	ARG	CA-CB-CG	-5.29	101.75	113.40
1	A	207	ALA	O-C-N	-5.25	114.29	122.70
1	B	9	LEU	N-CA-C	-5.25	96.82	111.00
1	B	89	VAL	CG1-CB-CG2	5.21	119.24	110.90
1	B	88	TRP	CG-CD1-NE1	-5.17	104.93	110.10
1	B	17	ILE	CA-CB-CG2	5.12	121.14	110.90
1	A	5	VAL	N-CA-C	-5.12	97.19	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	60	ARG	N-CA-CB	-5.01	101.59	110.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	100	TYR	Sidechain

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	1809	0	1755	19	0
1	B	1809	0	1755	30	0
2	A	13	0	14	0	0
2	B	13	0	14	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	186	0	0	3	0
5	B	174	0	0	4	0
All	All	4008	0	3538	46	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 6.

All (46) 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:55:ASN:HD22	1:B:57:VAL:H	1.30	0.79
1:B:129:MET:HG3	5:B:408:HOH:O	1.94	0.66
1:B:201:SER:HB2	1:B:206:PRO:HB3	1.79	0.64
1:B:55:ASN:ND2	1:B:58:ASP:H	1.95	0.64
1:B:55:ASN:HD21	1:B:58:ASP:H	1.43	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:ASN:ND2	1:B:131:ASN:H	1.97	0.62
1:A:55:ASN:HD22	1:A:57:VAL:H	1.52	0.58
1:B:187:VAL:HG23	1:B:188:VAL:HG23	1.88	0.56
1:A:157:THR:HB	5:A:332:HOH:O	2.05	0.56
1:A:43:GLN:HE21	1:A:46:LYS:HG3	1.71	0.55
1:B:54:TYR:HD2	5:B:389:HOH:O	1.88	0.55
1:B:43:GLN:HB3	1:B:46:LYS:HG3	1.90	0.53
1:A:201:SER:HB2	1:A:206:PRO:HB3	1.90	0.51
1:A:107:LEU:HD12	1:A:107:LEU:N	2.26	0.50
1:A:150:THR:HB	5:A:415:HOH:O	2.11	0.50
1:B:55:ASN:HD22	1:B:57:VAL:N	2.05	0.49
1:B:66:SER:HB3	1:B:72:SER:HB3	1.94	0.49
1:B:150:THR:HG21	5:B:279:HOH:O	2.12	0.49
1:A:117:SER:HA	1:A:186:ALA:HA	1.94	0.48
1:B:100:TYR:HB3	1:B:205:HIS:O	2.14	0.48
1:B:119:SER:O	1:B:122:GLU:HB2	2.13	0.48
1:B:13:PRO:HB3	1:B:20:PRO:O	2.14	0.48
1:A:112:THR:O	1:A:191:PHE:HA	2.13	0.48
1:B:86:PRO:HG2	1:B:89:VAL:HG23	1.96	0.47
1:B:228:ARG:HD3	5:B:256:HOH:O	2.15	0.47
1:B:115:LEU:O	1:B:123:THR:HA	2.14	0.47
1:A:131:ASN:H	1:B:124:ASN:ND2	2.12	0.47
1:A:32:VAL:HG13	1:A:233:PHE:CD2	2.51	0.45
1:B:32:VAL:HG13	1:B:233:PHE:CD2	2.51	0.45
1:B:90:ARG:HB3	1:B:174:LEU:HD12	1.98	0.45
1:B:3:THR:HG23	1:B:30:LYS:HD3	1.98	0.45
1:A:60:ARG:NH1	5:A:389:HOH:O	2.50	0.44
1:A:124:ASN:HD21	1:B:131:ASN:H	1.64	0.44
1:A:55:ASN:ND2	1:A:58:ASP:H	2.15	0.43
1:A:139:ASP:O	1:A:176:TYR:HB2	2.19	0.43
1:B:135:LYS:HB3	1:B:135:LYS:NZ	2.34	0.43
1:A:43:GLN:HB3	1:A:46:LYS:HG3	2.01	0.42
1:B:158:ARG:HH21	1:B:166:GLN:HG3	1.84	0.42
1:A:66:SER:HB3	1:A:72:SER:HB3	2.02	0.42
1:B:45:GLY:HA2	1:B:200:LYS:HB3	2.02	0.41
1:B:64:VAL:HG13	1:B:74:THR:HG22	2.02	0.41
1:B:55:ASN:ND2	1:B:57:VAL:HB	2.36	0.41
1:B:87:GLU:HG3	1:B:182:TRP:O	2.21	0.41
1:B:205:HIS:HA	1:B:206:PRO:HD2	1.82	0.40
1:A:89:VAL:HG22	1:A:215:SER:O	2.21	0.40
1:A:90:ARG:HB3	1:A:174:LEU:HD23	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	235/237 (99%)	224 (95%)	10 (4%)	1 (0%)	39	33
1	B	235/237 (99%)	223 (95%)	11 (5%)	1 (0%)	39	33
All	All	470/474 (99%)	447 (95%)	21 (4%)	2 (0%)	39	33

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	161	SER
1	B	150	THR

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	203/203 (100%)	185 (91%)	18 (9%)	12	7
1	B	203/203 (100%)	179 (88%)	24 (12%)	6	3
All	All	406/406 (100%)	364 (90%)	42 (10%)	9	5

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ILE

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Mol	Chain	Res	Type
1	A	32	VAL
1	A	69	ASN
1	A	81	LEU
1	A	89	VAL
1	A	101	LYS
1	A	120	THR
1	A	124	ASN
1	A	132	GLN
1	A	150	THR
1	A	151	ASP
1	A	158	ARG
1	A	162	ASN
1	A	184	SER
1	A	198	LEU
1	A	203	ASP
1	A	230	LEU
1	A	232	LEU
1	B	17	ILE
1	B	32	VAL
1	B	46	LYS
1	B	55	ASN
1	B	60	ARG
1	B	74	THR
1	B	81	LEU
1	B	89	VAL
1	B	119	SER
1	B	122	GLU
1	B	123	THR
1	B	124	ASN
1	B	126	LEU
1	B	129	MET
1	B	161	SER
1	B	174	LEU
1	B	181	ILE
1	B	183	GLU
1	B	198	LEU
1	B	203	ASP
1	B	228	ARG
1	B	229	LEU
1	B	230	LEU
1	B	232	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	55	ASN
1	A	104	ASN
1	A	124	ASN
1	A	131	ASN
1	A	162	ASN
1	B	43	GLN
1	B	55	ASN
1	B	104	ASN
1	B	124	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](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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	GYP	A	238	-	13,13,13	1.12	1 (7%)	18,18,18	4.72	2 (11%)
2	GYP	B	238	-	13,13,13	0.75	1 (7%)	18,18,18	0.89	1 (5%)

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	GYP	A	238	-	-	0/4/24/24	0/1/1/1
2	GYP	B	238	-	-	0/4/24/24	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	238	GYP	O1-C1	2.36	1.44	1.40
2	A	238	GYP	O1-C1	3.20	1.45	1.40

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	238	GYP	O4-C4-C3	-2.18	105.44	110.34
2	B	238	GYP	C7-O1-C1	2.25	116.98	113.29
2	A	238	GYP	C7-O1-C1	19.73	145.58	113.29

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 [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 [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, 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	237/237 (100%)	-0.49	10 (4%) 40 41	8, 17, 47, 70	0
1	B	237/237 (100%)	-0.49	6 (2%) 61 61	6, 15, 46, 76	0
All	All	474/474 (100%)	-0.49	16 (3%) 49 50	6, 16, 47, 76	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	119	SER	5.5
1	A	120	THR	5.5
1	A	1	ALA	4.3
1	B	121	HIS	3.7
1	B	118	ASN	3.7
1	A	162	ASN	3.5
1	B	120	THR	3.3
1	A	119	SER	3.2
1	A	118	ASN	3.0
1	B	162	ASN	2.9
1	A	161	SER	2.8
1	A	135	LYS	2.5
1	A	121	HIS	2.5
1	B	122	GLU	2.4
1	A	163	GLY	2.2
1	A	204	SER	2.1

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	GYP	A	238	13/13	0.97	0.08	0.93	18,21,24,25	0
2	GYP	B	238	13/13	0.98	0.08	0.26	14,17,17,17	0
3	MN	B	239	1/1	1.00	0.05	-1.01	14,14,14,14	0
3	MN	A	239	1/1	1.00	0.04	-1.60	13,13,13,13	0
4	CA	A	240	1/1	1.00	0.05	-2.26	13,13,13,13	0
4	CA	B	240	1/1	1.00	0.04	-2.39	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.