



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 07:56 PM BST

PDB ID : 5GUE
Title : Crystal structure of CotB2 (GGSP/Mg²⁺-Bound Form) from *Streptomyces melanosporofaciens*
Authors : Tomita, T.; Kim, S.-Y.; Ozaki, T.; Yoshida, A.; Kuzuyama, T.; Nishiyama, M.
Deposited on : 2016-08-28
Resolution : 1.80 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
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) : 2.13.1

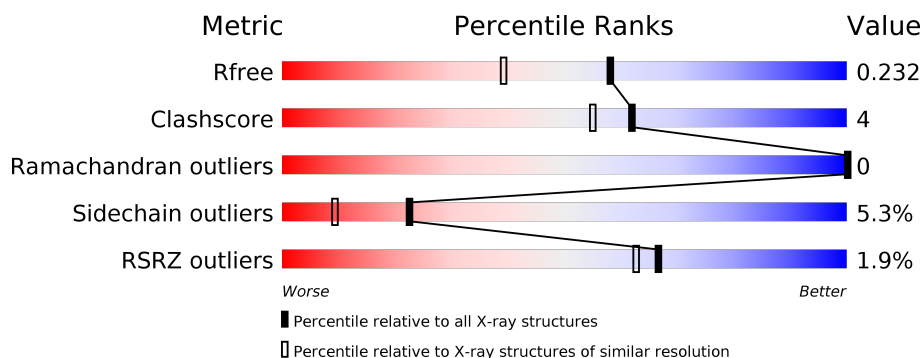
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 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	331	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>10%</div> <div>•</div> <div>14%</div> </div> </div>
1	B	331	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>9%</div> <div>•</div> <div>14%</div> </div> </div>
1	C	331	<div> <div>%</div> <div> <div></div> <div>76%</div> <div>8%</div> <div>•</div> <div>14%</div> </div> </div>
1	D	331	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>9%</div> <div>•</div> <div>14%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10326 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 Cyclooctat-9-en-7-ol synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	2	0
			2352	1499	398	438	17			
1	B	285	Total	C	N	O	S	0	0	0
			2336	1489	392	438	17			
1	C	285	Total	C	N	O	S	0	0	0
			2336	1489	392	438	17			
1	D	285	Total	C	N	O	S	0	0	0
			2336	1489	392	438	17			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	expression tag	UNP C9K1X5
A	-22	LYS	-	expression tag	UNP C9K1X5
A	-21	HIS	-	expression tag	UNP C9K1X5
A	-20	HIS	-	expression tag	UNP C9K1X5
A	-19	HIS	-	expression tag	UNP C9K1X5
A	-18	HIS	-	expression tag	UNP C9K1X5
A	-17	HIS	-	expression tag	UNP C9K1X5
A	-16	HIS	-	expression tag	UNP C9K1X5
A	-15	HIS	-	expression tag	UNP C9K1X5
A	-14	HIS	-	expression tag	UNP C9K1X5
A	-13	GLY	-	expression tag	UNP C9K1X5
A	-12	GLY	-	expression tag	UNP C9K1X5
A	-11	LEU	-	expression tag	UNP C9K1X5
A	-10	VAL	-	expression tag	UNP C9K1X5
A	-9	PRO	-	expression tag	UNP C9K1X5
A	-8	ARG	-	expression tag	UNP C9K1X5
A	-7	GLY	-	expression tag	UNP C9K1X5
A	-6	SER	-	expression tag	UNP C9K1X5
A	-5	HIS	-	expression tag	UNP C9K1X5
A	-4	GLY	-	expression tag	UNP C9K1X5
A	-3	GLY	-	expression tag	UNP C9K1X5

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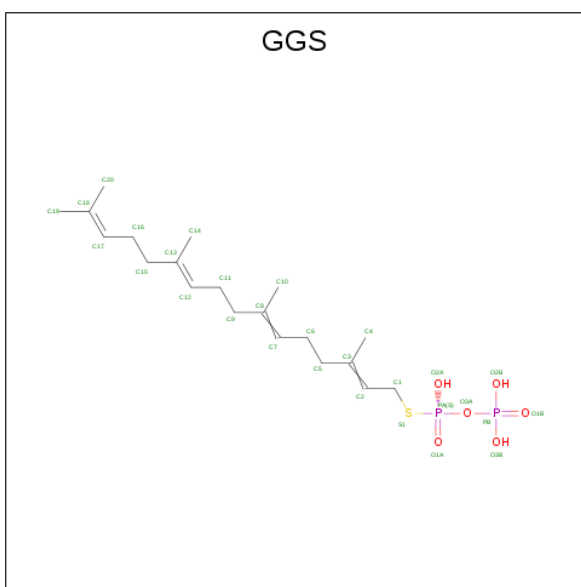
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP C9K1X5
A	-1	GLU	-	expression tag	UNP C9K1X5
A	0	PHE	-	expression tag	UNP C9K1X5
B	-23	MET	-	expression tag	UNP C9K1X5
B	-22	LYS	-	expression tag	UNP C9K1X5
B	-21	HIS	-	expression tag	UNP C9K1X5
B	-20	HIS	-	expression tag	UNP C9K1X5
B	-19	HIS	-	expression tag	UNP C9K1X5
B	-18	HIS	-	expression tag	UNP C9K1X5
B	-17	HIS	-	expression tag	UNP C9K1X5
B	-16	HIS	-	expression tag	UNP C9K1X5
B	-15	HIS	-	expression tag	UNP C9K1X5
B	-14	HIS	-	expression tag	UNP C9K1X5
B	-13	GLY	-	expression tag	UNP C9K1X5
B	-12	GLY	-	expression tag	UNP C9K1X5
B	-11	LEU	-	expression tag	UNP C9K1X5
B	-10	VAL	-	expression tag	UNP C9K1X5
B	-9	PRO	-	expression tag	UNP C9K1X5
B	-8	ARG	-	expression tag	UNP C9K1X5
B	-7	GLY	-	expression tag	UNP C9K1X5
B	-6	SER	-	expression tag	UNP C9K1X5
B	-5	HIS	-	expression tag	UNP C9K1X5
B	-4	GLY	-	expression tag	UNP C9K1X5
B	-3	GLY	-	expression tag	UNP C9K1X5
B	-2	SER	-	expression tag	UNP C9K1X5
B	-1	GLU	-	expression tag	UNP C9K1X5
B	0	PHE	-	expression tag	UNP C9K1X5
C	-23	MET	-	expression tag	UNP C9K1X5
C	-22	LYS	-	expression tag	UNP C9K1X5
C	-21	HIS	-	expression tag	UNP C9K1X5
C	-20	HIS	-	expression tag	UNP C9K1X5
C	-19	HIS	-	expression tag	UNP C9K1X5
C	-18	HIS	-	expression tag	UNP C9K1X5
C	-17	HIS	-	expression tag	UNP C9K1X5
C	-16	HIS	-	expression tag	UNP C9K1X5
C	-15	HIS	-	expression tag	UNP C9K1X5
C	-14	HIS	-	expression tag	UNP C9K1X5
C	-13	GLY	-	expression tag	UNP C9K1X5
C	-12	GLY	-	expression tag	UNP C9K1X5
C	-11	LEU	-	expression tag	UNP C9K1X5
C	-10	VAL	-	expression tag	UNP C9K1X5
C	-9	PRO	-	expression tag	UNP C9K1X5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-8	ARG	-	expression tag	UNP C9K1X5
C	-7	GLY	-	expression tag	UNP C9K1X5
C	-6	SER	-	expression tag	UNP C9K1X5
C	-5	HIS	-	expression tag	UNP C9K1X5
C	-4	GLY	-	expression tag	UNP C9K1X5
C	-3	GLY	-	expression tag	UNP C9K1X5
C	-2	SER	-	expression tag	UNP C9K1X5
C	-1	GLU	-	expression tag	UNP C9K1X5
C	0	PHE	-	expression tag	UNP C9K1X5
D	-23	MET	-	expression tag	UNP C9K1X5
D	-22	LYS	-	expression tag	UNP C9K1X5
D	-21	HIS	-	expression tag	UNP C9K1X5
D	-20	HIS	-	expression tag	UNP C9K1X5
D	-19	HIS	-	expression tag	UNP C9K1X5
D	-18	HIS	-	expression tag	UNP C9K1X5
D	-17	HIS	-	expression tag	UNP C9K1X5
D	-16	HIS	-	expression tag	UNP C9K1X5
D	-15	HIS	-	expression tag	UNP C9K1X5
D	-14	HIS	-	expression tag	UNP C9K1X5
D	-13	GLY	-	expression tag	UNP C9K1X5
D	-12	GLY	-	expression tag	UNP C9K1X5
D	-11	LEU	-	expression tag	UNP C9K1X5
D	-10	VAL	-	expression tag	UNP C9K1X5
D	-9	PRO	-	expression tag	UNP C9K1X5
D	-8	ARG	-	expression tag	UNP C9K1X5
D	-7	GLY	-	expression tag	UNP C9K1X5
D	-6	SER	-	expression tag	UNP C9K1X5
D	-5	HIS	-	expression tag	UNP C9K1X5
D	-4	GLY	-	expression tag	UNP C9K1X5
D	-3	GLY	-	expression tag	UNP C9K1X5
D	-2	SER	-	expression tag	UNP C9K1X5
D	-1	GLU	-	expression tag	UNP C9K1X5
D	0	PHE	-	expression tag	UNP C9K1X5

- Molecule 2 is phosphonoxy-[(10E)-3,7,11,15-tetramethylhexadeca-2,6,10,14-tetraenyl]sulfa nyl-phosphinic acid (three-letter code: GGS) (formula: C₂₀H₃₆O₆P₂S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 29	C 20	O 6	P 2	S 1	0	0
2	B	1	Total 29	C 20	O 6	P 2	S 1	0	0
2	C	1	Total 29	C 20	O 6	P 2	S 1	0	0
2	D	1	Total 29	C 20	O 6	P 2	S 1	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Mg 1 1	0	0
3	A	1	Total Mg 1 1	0	0
3	D	1	Total Mg 1 1	0	0
3	C	1	Total Mg 1 1	0	0

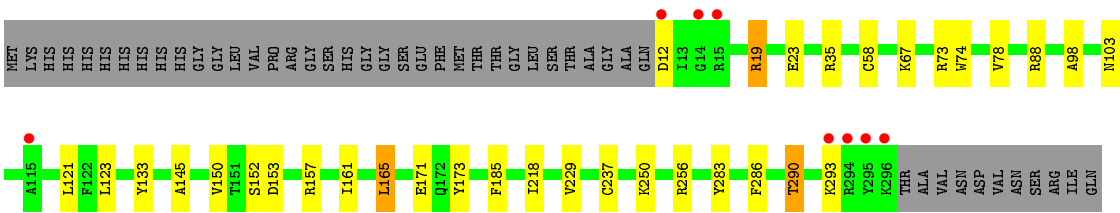
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	212	Total O 212 212	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	197	Total 197	O 197	0	0
4	C	217	Total 217	O 217	0	0
4	D	220	Total 220	O 220	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.70Å 100.37Å 108.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.15 – 1.80 34.15 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.1 (34.15-1.80) 99.0 (34.15-1.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 1.81Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.191 , 0.231 0.191 , 0.232	Depositor DCC
R_{free} test set	5798 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	26.0	Xtriage
Anisotropy	0.099	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 32.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.459 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10326	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 82.32 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.8711e-07. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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: MG, GGS

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	0.68	0/2416	0.71	1/3273 (0.0%)
1	B	0.69	0/2394	0.71	1/3245 (0.0%)
1	C	0.67	0/2394	0.71	0/3245
1	D	0.68	0/2394	0.69	0/3245
All	All	0.68	0/9598	0.71	2/13008 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	183	VAL	CB-CA-C	-7.37	97.40	111.40
1	B	183	VAL	CB-CA-C	-7.21	97.69	111.40

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	2352	0	2277	17	0
1	B	2336	0	2251	24	0
1	C	2336	0	2251	19	0
1	D	2336	0	2251	14	0
2	A	29	0	33	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	29	0	33	8	0
2	C	29	0	33	4	0
2	D	29	0	33	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	212	0	0	6	0
4	B	197	0	0	1	0
4	C	217	0	0	5	0
4	D	220	0	0	3	0
All	All	10326	0	9162	81	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 4.

All (81) 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:293:LYS:H	1:A:293:LYS:HD2	1.33	0.93
2:A:401:GGs:H11	2:A:401:GGs:H6	1.60	0.82
2:C:401:GGs:H11	2:C:401:GGs:H6	1.61	0.82
1:C:235:THR:HG23	4:C:606:HOH:O	1.79	0.81
1:B:178:VAL:HA	1:B:213:THR:HG21	1.63	0.80
1:A:62:VAL:HG13	1:A:67:LYS:HB3	1.64	0.79
2:D:401:GGs:H11	2:D:401:GGs:H6	1.73	0.70
1:B:183:VAL:HG21	1:B:213:THR:HG22	1.71	0.70
1:D:88:ARG:HD2	4:D:564:HOH:O	1.93	0.69
1:A:19:ARG:NH2	1:A:23:GLU:OE1	2.32	0.63
1:C:19:ARG:NH2	1:C:23:GLU:OE1	2.33	0.62
1:A:235:THR:HG23	4:A:566:HOH:O	2.00	0.61
1:B:281:LEU:HD11	2:B:401:GGs:H19A	1.82	0.61
1:B:183:VAL:HG13	1:B:216:LEU:CD1	2.30	0.61
1:C:161:ILE:HG22	1:C:165:LEU:HD22	1.88	0.56
2:B:401:GGs:H11	2:B:401:GGs:H6	1.87	0.56
1:D:19:ARG:NH2	1:D:23:GLU:OE1	2.39	0.56
1:A:102:LEU:O	1:A:106:THR:HG23	2.04	0.56
1:D:103:ASN:OD1	2:D:401:GGs:H9	2.06	0.56
1:A:152:SER:HB2	4:A:681:HOH:O	2.07	0.55
1:D:286:PHE:O	1:D:290:THR:HB	2.06	0.55
1:A:83:TRP:O	4:A:501:HOH:O	2.18	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:209:ALA:O	1:B:213:THR:CG2	2.56	0.54
1:D:161:ILE:HG22	1:D:165:LEU:HD22	1.90	0.53
1:C:91:LYS:NZ	4:C:504:HOH:O	2.36	0.53
1:C:80:VAL:HG21	2:C:401:GGG:H14B	1.91	0.53
1:B:103:ASN:OD1	2:B:401:GGG:H9	2.09	0.52
1:B:183:VAL:CG1	1:B:216:LEU:HD12	2.40	0.52
1:A:98:ALA:HB2	1:A:133:TYR:CB	2.40	0.52
1:C:181:ILE:CG2	1:C:183:VAL:HG23	2.40	0.51
1:B:209:ALA:O	1:B:213:THR:HG23	2.11	0.51
1:C:293:LYS:H	1:C:293:LYS:HD2	1.75	0.51
1:B:98:ALA:HB2	1:B:133:TYR:CB	2.41	0.50
1:C:67:LYS:HE3	1:C:112:MET:HG2	1.93	0.50
1:A:103:ASN:OD1	2:A:401:GGG:H9	2.12	0.50
1:C:62:VAL:HG13	1:C:67:LYS:HB2	1.92	0.50
1:C:98:ALA:HB2	1:C:133:TYR:CB	2.42	0.49
1:B:183:VAL:CG1	1:B:216:LEU:CD1	2.92	0.48
1:B:183:VAL:HG13	1:B:216:LEU:HD12	1.96	0.47
2:B:401:GGG:C18	2:B:401:GGG:H12	2.44	0.47
1:B:281:LEU:HD11	2:B:401:GGG:C19	2.43	0.47
1:C:43:LYS:HE2	1:C:82:LEU:HD23	1.97	0.47
1:C:173:TYR:CE2	1:C:237:CYS:HB3	2.50	0.47
1:B:113:ASP:HB3	1:B:116:LEU:HD12	1.97	0.46
1:A:173:TYR:CE2	1:A:237:CYS:HB3	2.50	0.46
1:B:181:ILE:CG2	1:B:183:VAL:HG22	2.45	0.46
1:B:34:ASP:OD2	1:C:34:ASP:OD2	2.33	0.45
1:A:106:THR:HG21	1:A:185:PHE:HE2	1.80	0.45
1:B:161:ILE:HG22	1:B:165:LEU:HD22	1.99	0.45
1:B:119:PHE:HE1	1:B:150:VAL:HG13	1.82	0.45
1:C:41:PRO:HD3	1:C:283:TYR:CE2	2.52	0.44
1:A:206:SER:HA	4:A:632:HOH:O	2.17	0.44
1:A:24:GLU:OE2	1:A:250:LYS:HE3	2.17	0.44
2:B:401:GGG:C11	2:B:401:GGG:H6	2.48	0.44
2:B:401:GGG:C17	2:B:401:GGG:H12	2.48	0.44
2:B:401:GGG:C6	2:B:401:GGG:H11	2.47	0.43
1:D:152:SER:HA	4:D:679:HOH:O	2.18	0.43
1:B:177:ARG:O	1:B:181:ILE:HG22	2.17	0.43
1:A:292:ASN:O	1:A:296:LYS:HE2	2.18	0.43
4:C:565:HOH:O	1:D:161:ILE:HD11	2.18	0.43
1:D:58:CYS:HB3	1:D:74:TRP:CH2	2.54	0.43
1:A:281:LEU:HD11	2:A:401:GGG:H19A	2.02	0.42
1:C:161:ILE:HD11	4:D:554:HOH:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:153:ASP:O	1:D:157:ARG:HG3	2.20	0.42
1:A:161:ILE:HD11	4:B:536:HOH:O	2.19	0.42
4:A:555:HOH:O	1:B:161:ILE:HD11	2.19	0.42
1:C:186:TRP:HB2	2:C:401:GGG:H10	2.01	0.42
1:D:98:ALA:HB2	1:D:133:TYR:CB	2.49	0.42
1:D:218:ILE:HD13	1:D:256:ARG:HB3	2.02	0.41
1:B:173:TYR:CE2	1:B:237:CYS:HB3	2.55	0.41
1:C:206:SER:HA	4:C:626:HOH:O	2.20	0.41
1:B:209:ALA:O	1:B:213:THR:HG22	2.20	0.41
1:D:173:TYR:CE2	1:D:237:CYS:HB3	2.55	0.41
1:D:58:CYS:HB3	1:D:74:TRP:CZ2	2.55	0.41
1:C:153:ASP:OD1	4:C:501:HOH:O	2.22	0.41
1:B:145:ALA:HB1	1:B:185:PHE:HA	2.03	0.40
1:B:13:ILE:O	1:B:229:VAL:HG13	2.22	0.40
1:C:182:GLY:O	2:C:401:GGG:H10B	2.21	0.40
1:D:145:ALA:HB1	1:D:185:PHE:HA	2.04	0.40
1:A:235:THR:HG22	4:A:664:HOH:O	2.20	0.40
1:B:191:TYR:CG	1:B:192:PRO:HD3	2.55	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	285/331 (86%)	284 (100%)	1 (0%)	0	100	100
1	B	283/331 (86%)	281 (99%)	2 (1%)	0	100	100
1	C	283/331 (86%)	279 (99%)	4 (1%)	0	100	100
1	D	283/331 (86%)	281 (99%)	2 (1%)	0	100	100
All	All	1134/1324 (86%)	1125 (99%)	9 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	258/292 (88%)	243 (94%)	15 (6%)	20	7
1	B	256/292 (88%)	243 (95%)	13 (5%)	24	10
1	C	256/292 (88%)	245 (96%)	11 (4%)	29	14
1	D	256/292 (88%)	240 (94%)	16 (6%)	18	6
All	All	1026/1168 (88%)	971 (95%)	55 (5%)	22	9

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27[A]	ARG
1	A	27[B]	ARG
1	A	67	LYS
1	A	68	THR
1	A	73	ARG
1	A	78	VAL
1	A	121	LEU
1	A	126	LEU
1	A	150	VAL
1	A	155	MET
1	A	158	ASP
1	A	171	GLU
1	A	183	VAL
1	A	283	TYR
1	A	293	LYS
1	B	67	LYS
1	B	68	THR
1	B	78	VAL
1	B	121	LEU
1	B	123	LEU
1	B	126	LEU
1	B	150	VAL
1	B	165	LEU
1	B	183	VAL

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Mol	Chain	Res	Type
1	B	213	THR
1	B	229	VAL
1	B	247	THR
1	B	283	TYR
1	C	19	ARG
1	C	44	VAL
1	C	82	LEU
1	C	121	LEU
1	C	123	LEU
1	C	126	LEU
1	C	165	LEU
1	C	171	GLU
1	C	247	THR
1	C	283	TYR
1	C	293	LYS
1	D	12	ASP
1	D	19	ARG
1	D	35	ARG
1	D	67	LYS
1	D	73	ARG
1	D	78	VAL
1	D	121	LEU
1	D	123	LEU
1	D	150	VAL
1	D	165	LEU
1	D	171	GLU
1	D	229	VAL
1	D	250	LYS
1	D	283	TYR
1	D	290	THR
1	D	293	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	HIS
1	A	172	GLN
1	A	233	GLN
1	B	172	GLN
1	C	172	GLN
1	C	233	GLN
1	D	172	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	GGS	C	401	3	24,28,28	2.11	9 (37%)	29,37,37	1.46	3 (10%)
2	GGS	D	401	3	24,28,28	2.14	10 (41%)	29,37,37	1.55	7 (24%)
2	GGS	A	401	3	24,28,28	2.17	10 (41%)	29,37,37	1.53	5 (17%)
2	GGS	B	401	3	24,28,28	2.13	9 (37%)	29,37,37	1.40	5 (17%)

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	GGS	C	401	3	-	4/25/31/31	-
2	GGS	D	401	3	-	6/25/31/31	-
2	GGS	A	401	3	-	7/25/31/31	-
2	GGS	B	401	3	-	8/25/31/31	-

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	GGs	PB-O1B	3.99	1.63	1.50
2	B	401	GGs	PB-O1B	3.76	1.62	1.50
2	A	401	GGs	C16-C17	-3.75	1.38	1.50
2	D	401	GGs	C16-C17	-3.71	1.38	1.50
2	D	401	GGs	PB-O1B	3.66	1.62	1.50
2	C	401	GGs	PB-O1B	3.65	1.62	1.50
2	B	401	GGs	C16-C17	-3.50	1.39	1.50
2	A	401	GGs	C6-C7	-3.33	1.39	1.50
2	B	401	GGs	C11-C12	-3.32	1.39	1.50
2	B	401	GGs	PA-O2A	-3.31	1.48	1.56
2	B	401	GGs	C6-C7	-3.29	1.39	1.50
2	A	401	GGs	C11-C12	-3.29	1.39	1.50
2	C	401	GGs	C6-C7	-3.27	1.39	1.50
2	C	401	GGs	C16-C17	-3.25	1.39	1.50
2	D	401	GGs	C6-C7	-3.25	1.39	1.50
2	C	401	GGs	C11-C12	-3.24	1.39	1.50
2	D	401	GGs	C11-C12	-3.17	1.40	1.50
2	D	401	GGs	C7-C8	3.12	1.40	1.33
2	C	401	GGs	C7-C8	3.12	1.40	1.33
2	C	401	GGs	PA-O2A	-3.10	1.48	1.56
2	A	401	GGs	C12-C13	3.09	1.40	1.33
2	D	401	GGs	PA-O2A	-3.09	1.48	1.56
2	D	401	GGs	C12-C13	3.06	1.40	1.33
2	C	401	GGs	C2-C3	3.05	1.40	1.33
2	B	401	GGs	C12-C13	3.03	1.40	1.33
2	A	401	GGs	PA-O2A	-3.00	1.49	1.56
2	C	401	GGs	C12-C13	2.99	1.40	1.33
2	A	401	GGs	C7-C8	2.98	1.40	1.33
2	B	401	GGs	C7-C8	2.97	1.40	1.33
2	D	401	GGs	C2-C3	2.96	1.40	1.33
2	A	401	GGs	C2-C3	2.95	1.40	1.33
2	B	401	GGs	C2-C3	2.86	1.39	1.33
2	C	401	GGs	C17-C18	2.66	1.40	1.32
2	B	401	GGs	C17-C18	2.32	1.39	1.32
2	A	401	GGs	C17-C18	2.26	1.38	1.32
2	D	401	GGs	C17-C18	2.21	1.38	1.32
2	A	401	GGs	C9-C8	2.02	1.55	1.51
2	D	401	GGs	PB-O2B	2.00	1.62	1.54

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	GGs	O3B-PB-O3A	3.30	115.71	104.64
2	A	401	GGs	O3B-PB-O3A	3.27	115.59	104.64
2	C	401	GGs	C9-C8-C7	3.12	127.43	121.12
2	D	401	GGs	C20-C18-C19	3.08	121.41	114.60
2	D	401	GGs	C9-C8-C7	3.05	127.28	121.12
2	C	401	GGs	C10-C8-C7	-3.02	115.94	123.68
2	A	401	GGs	C10-C8-C7	-3.00	115.99	123.68
2	B	401	GGs	O3B-PB-O3A	2.97	114.58	104.64
2	A	401	GGs	C20-C18-C19	2.89	120.98	114.60
2	A	401	GGs	C9-C8-C7	2.88	126.94	121.12
2	D	401	GGs	O3B-PB-O3A	2.86	114.22	104.64
2	D	401	GGs	C10-C8-C7	-2.61	116.99	123.68
2	B	401	GGs	C10-C8-C7	-2.60	117.00	123.68
2	D	401	GGs	C19-C18-C17	-2.54	115.31	122.65
2	D	401	GGs	C4-C3-C2	-2.53	117.19	123.68
2	B	401	GGs	C4-C3-C5	2.51	119.48	115.27
2	A	401	GGs	C19-C18-C17	-2.49	115.45	122.65
2	B	401	GGs	C9-C8-C7	2.47	126.11	121.12
2	B	401	GGs	C4-C3-C2	-2.45	117.40	123.68
2	D	401	GGs	C4-C3-C5	2.28	119.10	115.27

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	GGs	C12-C11-C9-C8
2	C	401	GGs	C3-C5-C6-C7
2	A	401	GGs	C3-C5-C6-C7
2	D	401	GGs	C10-C8-C9-C11
2	B	401	GGs	C10-C8-C9-C11
2	A	401	GGs	C7-C8-C9-C11
2	A	401	GGs	C10-C8-C9-C11
2	C	401	GGs	C7-C8-C9-C11
2	D	401	GGs	C7-C8-C9-C11
2	B	401	GGs	C7-C8-C9-C11
2	C	401	GGs	C10-C8-C9-C11
2	D	401	GGs	C13-C15-C16-C17
2	B	401	GGs	C3-C5-C6-C7
2	A	401	GGs	PA-O3A-PB-O1B
2	D	401	GGs	C3-C5-C6-C7
2	B	401	GGs	C13-C15-C16-C17
2	B	401	GGs	C14-C13-C15-C16
2	B	401	GGs	C12-C13-C15-C16

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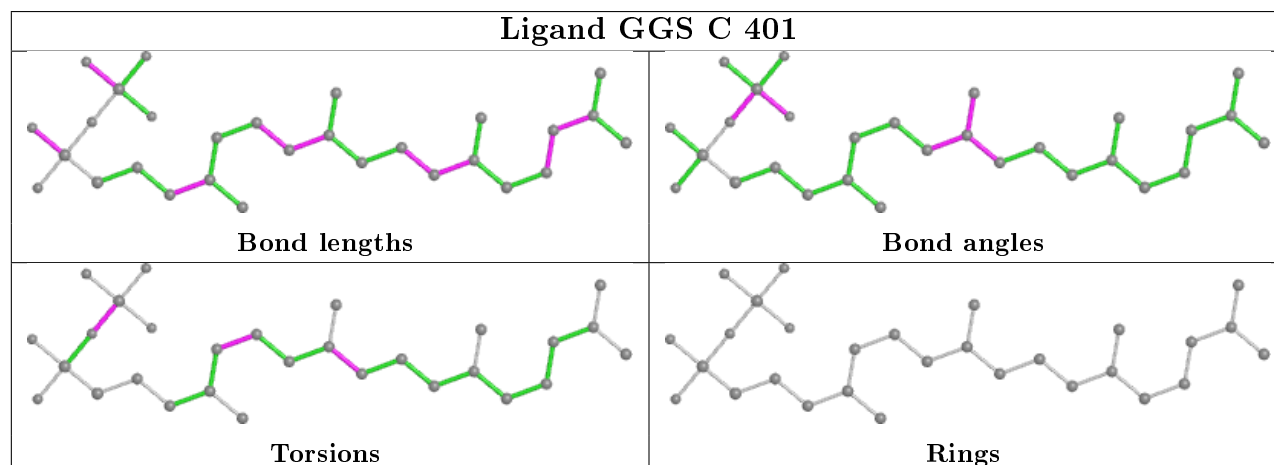
Mol	Chain	Res	Type	Atoms
2	A	401	GGG	C14-C13-C15-C16
2	D	401	GGG	C14-C13-C15-C16
2	C	401	GGG	PA-O3A-PB-O1B
2	A	401	GGG	C12-C11-C9-C8
2	D	401	GGG	C12-C13-C15-C16
2	A	401	GGG	C12-C13-C15-C16
2	B	401	GGG	C9-C11-C12-C13

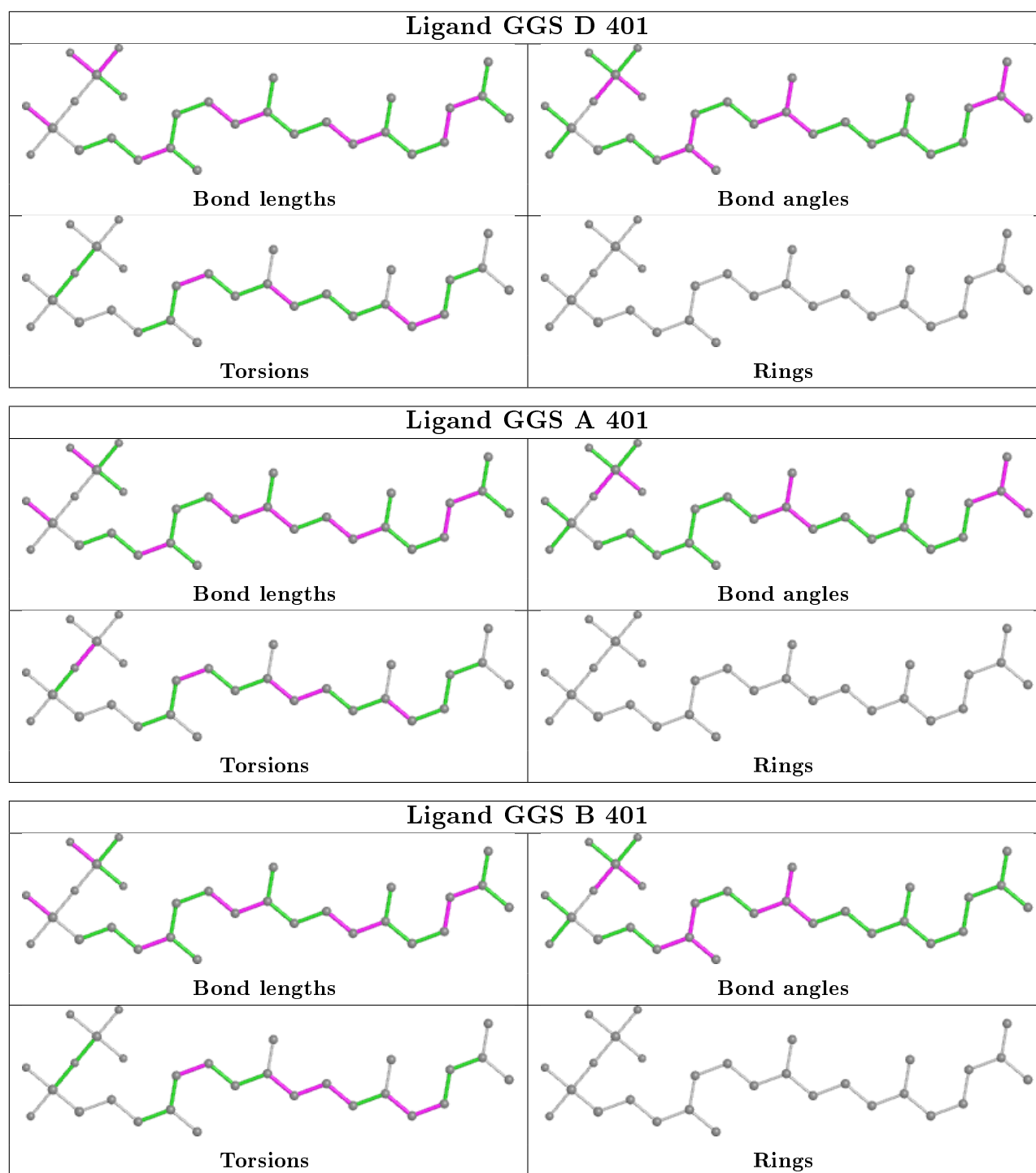
There are no ring outliers.

4 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	GGG	4	0
2	D	401	GGG	2	0
2	A	401	GGG	3	0
2	B	401	GGG	8	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	285/331 (86%)	-0.32	5 (1%) 68 64	18, 28, 45, 58	0
1	B	285/331 (86%)	-0.31	5 (1%) 68 64	18, 29, 45, 61	0
1	C	285/331 (86%)	-0.29	4 (1%) 75 72	18, 28, 46, 60	0
1	D	285/331 (86%)	-0.25	8 (2%) 53 47	18, 29, 45, 62	0
All	All	1140/1324 (86%)	-0.29	22 (1%) 66 63	18, 29, 46, 62	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	295	TYR	5.7
1	A	295	TYR	5.3
1	D	12	ASP	4.8
1	A	293	LYS	4.2
1	B	295	TYR	3.9
1	D	15	ARG	3.8
1	C	15	ARG	3.8
1	A	12	ASP	3.7
1	B	293	LYS	3.6
1	B	15	ARG	3.6
1	C	12	ASP	3.5
1	D	295	TYR	3.3
1	C	293	LYS	3.1
1	B	296	LYS	3.1
1	A	156	PHE	2.9
1	D	115	ALA	2.9
1	B	12	ASP	2.9
1	D	294	ARG	2.7
1	D	293	LYS	2.7
1	D	14	GLY	2.5
1	A	15	ARG	2.4
1	D	296	LYS	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 monosaccharides in this entry.

6.4 Ligands [i](#)

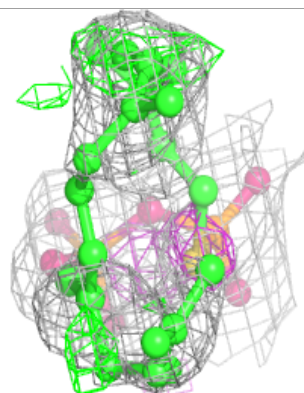
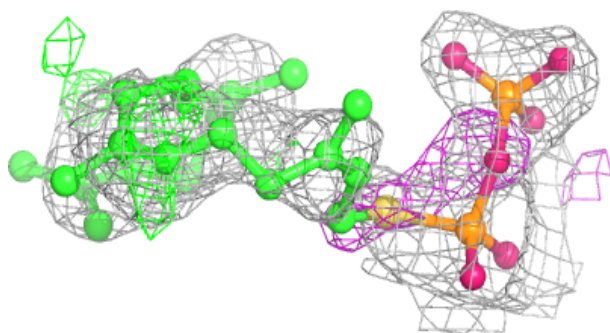
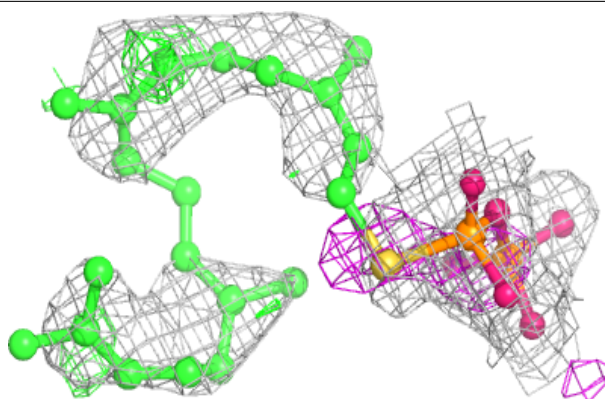
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	GGs	B	401	29/29	0.80	0.21	62,68,69,69	0
2	GGs	D	401	29/29	0.81	0.21	51,61,68,70	0
2	GGs	C	401	29/29	0.86	0.16	38,54,60,62	0
2	GGs	A	401	29/29	0.86	0.19	50,57,61,62	0
3	MG	B	402	1/1	0.95	0.12	39,39,39,39	0
3	MG	D	402	1/1	0.96	0.08	35,35,35,35	0
3	MG	A	402	1/1	0.97	0.04	34,34,34,34	0
3	MG	C	402	1/1	0.99	0.14	35,35,35,35	0

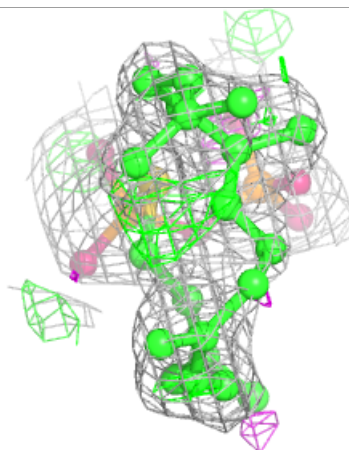
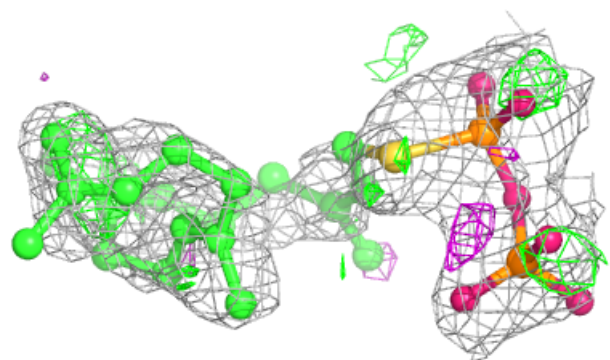
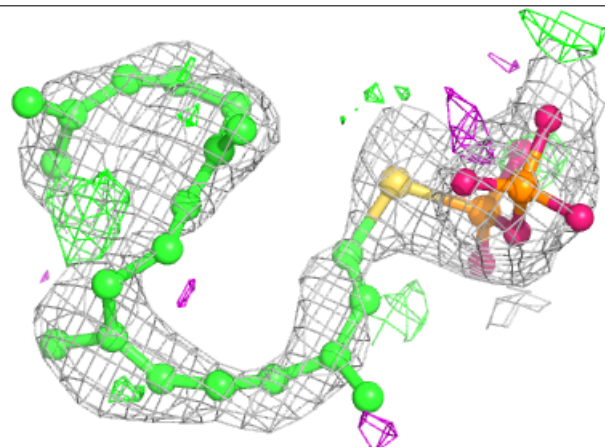
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around GGS B 401:

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

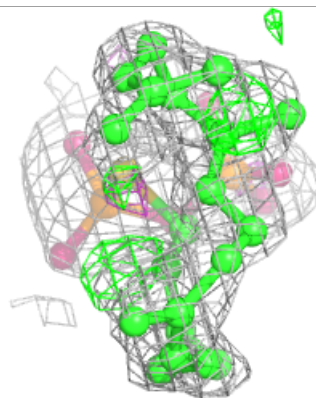
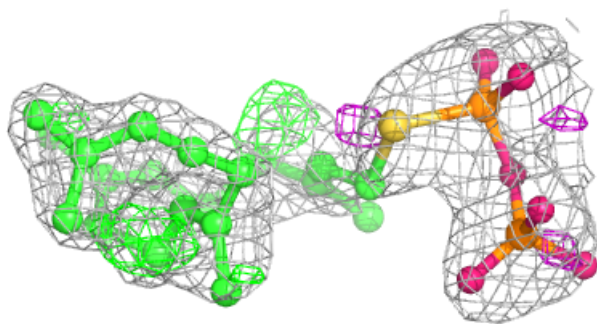
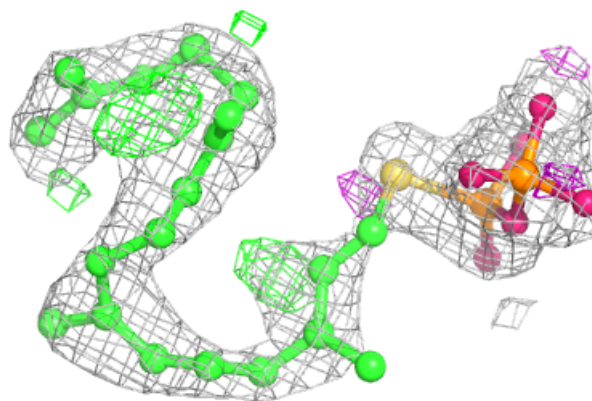
**Electron density around GGS D 401:**

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

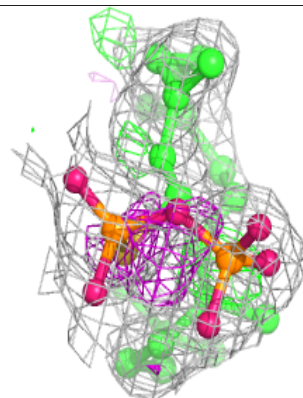
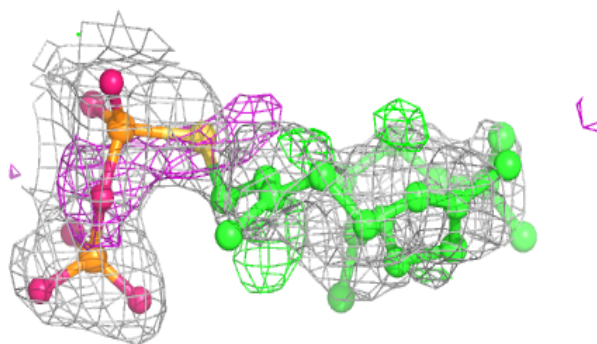
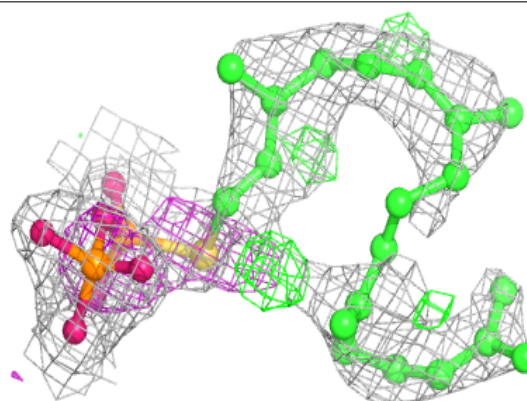


Electron density around GGS C 401:

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

**Electron density around GGS A 401:**

$2mF_o-DF_c$ (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.