



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 11:26 PM UTC

PDB ID : 3AE0 / pdb_00003ae0
Title : Crystal structure of the C(30) carotenoid dehydrosqualene synthase from *Staphylococcus aureus* complexed with geranylgeranyl thiopyrophosphate
Authors : Liu, C.I.; Jeng, W.Y.; Wang, A.H.J.; Oldfield, E.
Deposited on : 2010-01-31
Resolution : 2.37 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

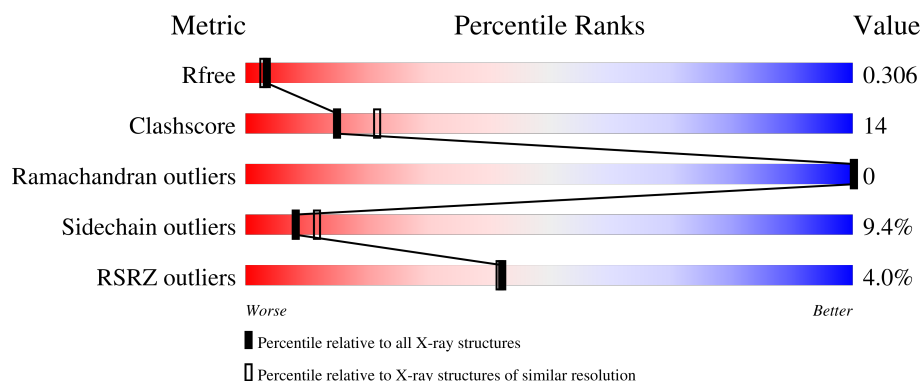
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

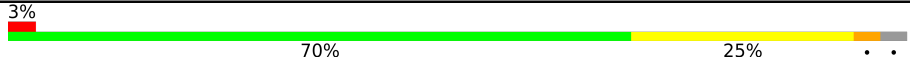
The reported resolution of this entry is 2.37 Å.

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}	180053	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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 of 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	293	
1	B	293	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GGs	A	1001	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4964 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.

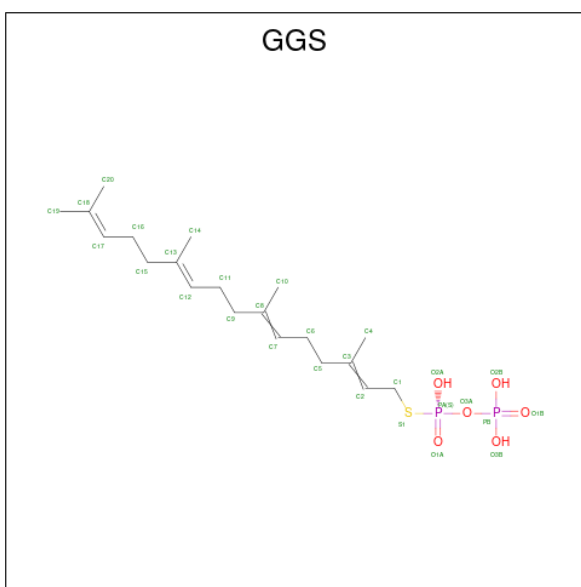
- Molecule 1 is a protein called Dehydrosqualene synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	284	Total	C	N	O	S	0	0	0
			2386	1530	400	444	12			
1	B	284	Total	C	N	O	S	0	0	0
			2386	1530	400	444	12			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	ALA	-	expression tag	UNP A9JQL9
A	-4	ALA	-	expression tag	UNP A9JQL9
A	-3	ALA	-	expression tag	UNP A9JQL9
A	-2	ALA	-	expression tag	UNP A9JQL9
A	-1	ALA	-	expression tag	UNP A9JQL9
A	0	ALA	-	expression tag	UNP A9JQL9
A	26	ALA	PHE	engineered mutation	UNP A9JQL9
B	-5	ALA	-	expression tag	UNP A9JQL9
B	-4	ALA	-	expression tag	UNP A9JQL9
B	-3	ALA	-	expression tag	UNP A9JQL9
B	-2	ALA	-	expression tag	UNP A9JQL9
B	-1	ALA	-	expression tag	UNP A9JQL9
B	0	ALA	-	expression tag	UNP A9JQL9
B	26	ALA	PHE	engineered mutation	UNP A9JQL9

- Molecule 2 is phosphonooxy-[(10E)-3,7,11,15-tetramethylhexadeca-2,6,10,14-tetraenyl]sulfonyl-phosphinic acid (CCD ID: GGS) (formula: C₂₀H₃₆O₆P₂S).



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

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

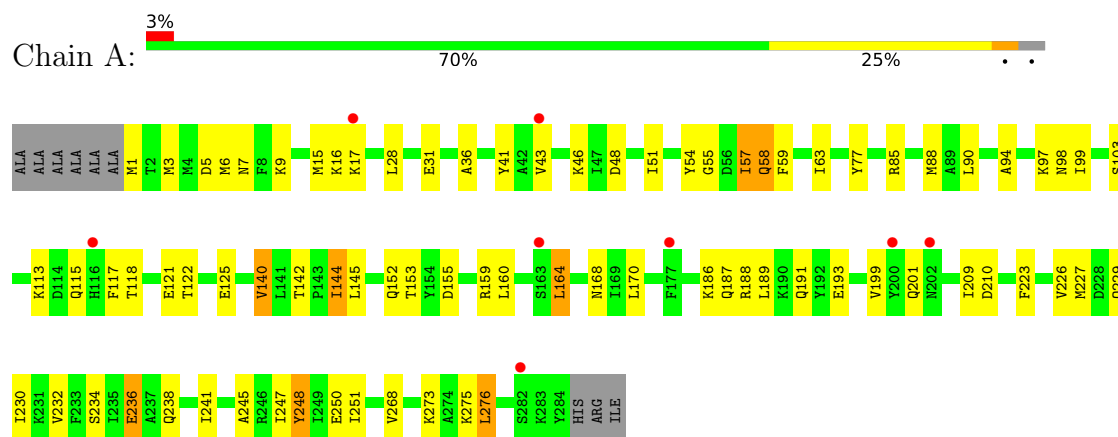
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	30	Total	O	0	0
			30	30		
4	B	40	Total	O	0	0
			40	40		

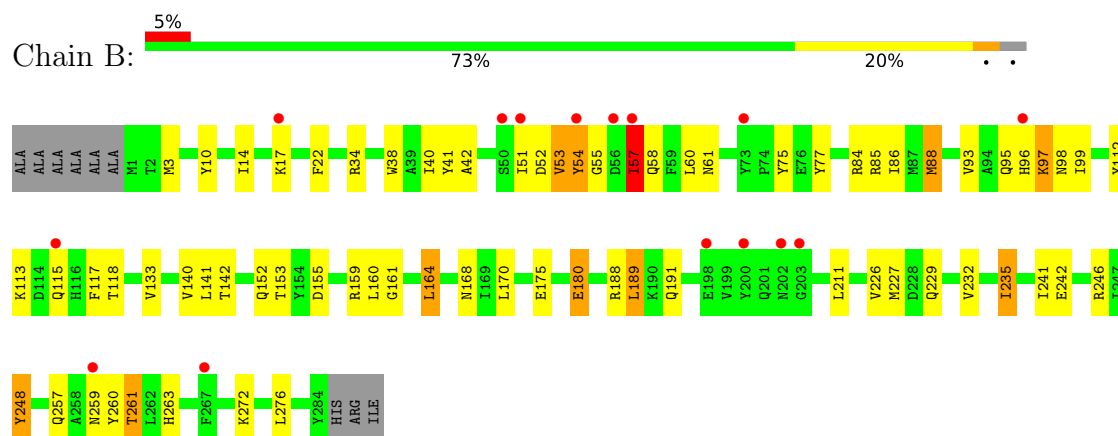
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Dehydrosqualene synthase



• Molecule 1: Dehydrosqualene synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	80.12Å 80.12Å 183.62Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.37 30.00 – 2.37	Depositor EDS
% Data completeness (in resolution range)	99.0 (30.00-2.37) 99.0 (30.00-2.37)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.12 (at 2.36Å)	Xtriage
Refinement program	CNS, REFMAC 5.5.0072	Depositor
R, R_{free}	0.219 , 0.281 0.260 , 0.306	Depositor DCC
R_{free} test set	1428 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	50.4	Xtriage
Anisotropy	0.347	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 32.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.037 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4964	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 90.11 % 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 2.6219e-08. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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	0.45	0/2441	0.96	0/3292
1	B	0.43	0/2441	0.97	1/3292 (0.0%)
All	All	0.44	0/4882	0.97	1/6584 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	57	ILE	N-CA-C	5.21	115.98	110.72

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	2386	0	2317	49	0
1	B	2386	0	2317	53	0
2	A	58	0	66	30	0
2	B	58	0	66	28	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
4	A	30	0	0	2	0
4	B	40	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4964	0	4766	132	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (132) 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)
2:B:1003:GGS:H19B	2:B:1004:GGS:C19	1.85	1.06
2:B:1003:GGS:C19	2:B:1004:GGS:H19B	1.87	1.04
2:B:1003:GGS:H19B	2:B:1004:GGS:H19B	1.41	1.01
2:A:1001:GGS:H19	2:A:1002:GGS:C18	1.94	0.97
1:B:152:GLN:HE21	1:B:229:GLN:HE21	1.22	0.87
1:A:6:MET:HG2	4:B:544:HOH:O	1.74	0.85
1:A:188:ARG:HH11	1:A:191:GLN:HE22	1.24	0.83
2:B:1003:GGS:C19	2:B:1004:GGS:C19	2.49	0.83
1:A:152:GLN:HE21	1:A:229:GLN:HE21	1.28	0.77
1:B:164:LEU:HB3	2:B:1003:GGS:C6	2.17	0.73
1:B:235:ILE:O	1:B:235:ILE:HD13	1.89	0.73
1:A:268:VAL:HG23	1:A:273:LYS:HE3	1.72	0.71
2:A:1001:GGS:H19A	2:A:1002:GGS:H19B	1.73	0.71
1:A:41:TYR:CD1	2:A:1002:GGS:H11	2.27	0.70
1:A:247:ILE:HG23	1:A:268:VAL:HG21	1.74	0.69
2:A:1001:GGS:C19	2:A:1002:GGS:H19B	2.22	0.69
1:B:175:GLU:HG3	4:B:531:HOH:O	1.93	0.69
1:B:164:LEU:HB3	2:B:1003:GGS:H6A	1.75	0.67
1:B:41:TYR:CE1	2:B:1004:GGS:H10	2.30	0.67
1:A:9:LYS:NZ	1:B:180:GLU:OE2	2.26	0.67
2:B:1004:GGS:H1	2:B:1004:GGS:O3B	1.94	0.67
1:B:141:LEU:HD11	2:B:1004:GGS:H14B	1.75	0.67
1:A:168:ASN:HB2	2:A:1001:GGS:H4B	1.76	0.67
2:B:1003:GGS:H19B	2:B:1004:GGS:H19	1.76	0.65
2:A:1001:GGS:H19	2:A:1002:GGS:C19	2.27	0.65
1:A:48:ASP:CG	2:A:1002:GGS:H2	2.22	0.64
2:A:1001:GGS:C19	2:A:1002:GGS:C19	2.76	0.64
2:A:1001:GGS:C2	2:A:1002:GGS:H4B	2.28	0.64
1:B:41:TYR:CD1	2:B:1004:GGS:H11	2.34	0.63
1:B:34:ARG:HD2	1:B:38:TRP:CZ2	2.34	0.62
1:B:53:VAL:HG22	1:B:54:TYR:HD1	1.65	0.62
2:B:1003:GGS:H19	2:B:1004:GGS:C18	2.29	0.62
1:A:155:ASP:O	1:A:159:ARG:HG3	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1004:GGS:H4	2:B:1004:GGS:C7	2.29	0.61
1:A:103:SER:HB2	1:A:140:VAL:HG23	1.81	0.61
1:B:142:THR:HG23	1:B:153:THR:HG22	1.82	0.61
2:B:1003:GGS:H1	2:B:1004:GGS:S1	2.42	0.60
1:A:188:ARG:NH1	1:A:191:GLN:HE22	1.96	0.60
1:B:112:TYR:C	1:B:112:TYR:CD2	2.80	0.59
1:B:113:LYS:C	1:B:115:GLN:H	2.11	0.59
1:B:77:TYR:OH	1:B:84:ARG:NH1	2.36	0.59
2:B:1003:GGS:C19	2:B:1004:GGS:C18	2.83	0.57
1:A:164:LEU:HB3	2:A:1001:GGS:H6A	1.87	0.56
2:B:1004:GGS:S1	2:B:1004:GGS:H4B	2.45	0.56
2:A:1001:GGS:H9A	2:A:1002:GGS:H12	1.87	0.56
1:B:57:ILE:HG13	1:B:57:ILE:O	2.07	0.55
1:B:42:ALA:HB1	1:B:86:ILE:HD12	1.88	0.54
1:A:226:VAL:HG12	1:A:227:MET:HE2	1.90	0.54
2:A:1001:GGS:C13	2:A:1001:GGS:H20B	2.37	0.54
1:A:15:MET:HE1	2:A:1002:GGS:H16	1.89	0.54
2:B:1004:GGS:H4	2:B:1004:GGS:H7	1.89	0.53
2:A:1001:GGS:H20B	2:A:1001:GGS:C12	2.38	0.53
2:A:1001:GGS:H19	2:A:1002:GGS:C20	2.39	0.53
1:B:168:ASN:ND2	2:B:1003:GGS:H2	2.24	0.53
1:A:115:GLN:C	1:A:117:PHE:H	2.17	0.52
2:B:1003:GGS:H10	2:B:1004:GGS:H9A	1.91	0.52
1:B:10:TYR:CE2	1:B:14:ILE:HD11	2.44	0.52
2:A:1001:GGS:H1	2:A:1002:GGS:H4B	1.90	0.52
1:B:164:LEU:HB3	2:B:1003:GGS:H6	1.90	0.52
1:A:236:GLU:H	1:A:236:GLU:CD	2.17	0.52
1:A:223:PHE:HE1	1:A:245:ALA:HB1	1.75	0.51
1:B:10:TYR:CZ	1:B:14:ILE:HD11	2.45	0.51
1:A:46:LYS:HE2	4:A:541:HOH:O	2.09	0.51
1:B:160:LEU:HD23	2:B:1003:GGS:H16A	1.92	0.51
1:A:57:ILE:HG13	1:A:58:GLN:N	2.22	0.51
1:A:223:PHE:CE1	1:A:245:ALA:HB1	2.46	0.51
1:A:51:ILE:O	1:A:55:GLY:HA2	2.11	0.50
1:B:58:GLN:HA	1:B:61:ASN:HD22	1.77	0.50
1:B:160:LEU:HD13	1:B:226:VAL:HG11	1.93	0.50
2:B:1003:GGS:H9A	2:B:1004:GGS:H12	1.95	0.49
1:A:187:GLN:NE2	4:A:557:HOH:O	2.45	0.49
1:A:250:GLU:HG3	1:A:276:LEU:HD21	1.96	0.48
1:A:43:VAL:HG21	1:A:90:LEU:HD22	1.96	0.48
1:A:41:TYR:HB2	2:A:1002:GGS:H14A	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1001:GGS:H2	2:A:1002:GGS:H4B	1.95	0.48
1:A:241:ILE:HG12	2:A:1001:GGS:H17	1.96	0.47
1:B:226:VAL:HG12	1:B:227:MET:HE2	1.95	0.47
1:A:59:PHE:O	1:A:63:ILE:HG13	2.13	0.47
1:A:122:THR:HG23	1:A:125:GLU:H	1.79	0.47
2:A:1002:GGS:C4	2:A:1002:GGS:H7	2.44	0.47
1:B:42:ALA:CB	1:B:86:ILE:HD12	2.45	0.47
1:B:152:GLN:HE21	1:B:229:GLN:NE2	2.02	0.47
1:A:41:TYR:CE1	2:A:1002:GGS:H11	2.50	0.47
1:B:188:ARG:HH11	1:B:191:GLN:HE22	1.63	0.46
1:A:77:TYR:CE1	1:A:88:MET:HE1	2.51	0.46
1:A:160:LEU:HG	2:A:1001:GGS:C14	2.46	0.46
1:A:51:ILE:O	1:A:55:GLY:N	2.48	0.46
1:B:160:LEU:HG	2:B:1003:GGS:H14B	1.97	0.45
1:B:242:GLU:OE2	1:B:246:ARG:NH1	2.49	0.45
1:B:188:ARG:HD3	1:B:191:GLN:NE2	2.32	0.45
1:B:75:TYR:CE1	1:B:95:GLN:HG2	2.52	0.45
1:B:40:ILE:HD12	1:B:140:VAL:CG1	2.48	0.44
1:B:235:ILE:HD13	1:B:235:ILE:C	2.42	0.44
1:B:53:VAL:HG22	1:B:54:TYR:CD1	2.49	0.44
1:B:112:TYR:C	1:B:112:TYR:HD2	2.23	0.44
1:B:51:ILE:O	1:B:52:ASP:C	2.61	0.44
1:A:7:ASN:ND2	1:A:85:ARG:HG2	2.33	0.43
1:B:96:HIS:C	1:B:97:LYS:HG2	2.43	0.43
1:B:14:ILE:HD12	1:B:86:ILE:HD11	2.00	0.43
1:B:259:ASN:C	1:B:260:TYR:CD2	2.97	0.43
1:A:142:THR:HG23	1:A:153:THR:HG22	2.00	0.43
1:B:115:GLN:C	1:B:117:PHE:H	2.27	0.43
1:A:94:ALA:HB2	1:A:99:ILE:HD12	2.01	0.43
2:A:1002:GGS:H16A	2:A:1002:GGS:H14	1.71	0.43
1:B:241:ILE:HG23	2:B:1003:GGS:H17	2.00	0.43
1:A:1:MET:HG3	1:A:5:ASP:CB	2.49	0.42
2:A:1001:GGS:C1	2:A:1002:GGS:H4B	2.49	0.42
1:B:51:ILE:O	1:B:55:GLY:HA2	2.19	0.42
1:B:84:ARG:O	1:B:88:MET:HB2	2.19	0.42
1:A:160:LEU:HD23	2:A:1001:GGS:H16A	2.01	0.42
1:A:234:SER:O	1:A:238:GLN:HG3	2.19	0.42
1:B:93:VAL:HG12	1:B:99:ILE:HD11	2.01	0.42
1:B:188:ARG:HD3	1:B:191:GLN:HE21	1.84	0.42
1:B:161:GLY:N	2:B:1003:GGS:H14A	2.34	0.42
1:A:36:ALA:HB1	1:A:144:ILE:HG12	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:GLU:HA	1:A:186:LYS:HB2	2.02	0.42
1:B:189:LEU:HD13	1:B:211:LEU:HD21	2.01	0.41
2:B:1003:GGS:C5	2:B:1003:GGS:S1	3.08	0.41
1:A:41:TYR:CE1	2:A:1002:GGS:H10	2.56	0.41
1:A:209:ILE:O	1:A:210:ASP:C	2.63	0.41
2:A:1001:GGS:S1	2:A:1001:GGS:C5	3.08	0.41
1:B:155:ASP:O	1:B:159:ARG:HG3	2.20	0.41
1:A:145:LEU:HD11	2:A:1001:GGS:H19A	2.03	0.41
1:A:168:ASN:HD21	2:A:1002:GGS:PA	2.44	0.41
1:A:188:ARG:HA	1:A:188:ARG:HD3	1.93	0.41
1:B:41:TYR:CE1	2:B:1004:GGS:H11	2.55	0.41
1:B:241:ILE:HG12	2:B:1003:GGS:H17	2.03	0.41
1:A:160:LEU:HG	2:A:1001:GGS:H14	2.03	0.40
1:B:261:THR:HG22	1:B:263:HIS:H	1.86	0.40
1:B:22:PHE:CZ	1:B:248:TYR:CE1	3.09	0.40
1:A:113:LYS:C	1:A:115:GLN:H	2.29	0.40
1:A:248:TYR:O	1:A:251:ILE:HB	2.22	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	282/293 (96%)	273 (97%)	9 (3%)	0	100	100
1	B	282/293 (96%)	270 (96%)	12 (4%)	0	100	100
All	All	564/586 (96%)	543 (96%)	21 (4%)	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	256/259 (99%)	231 (90%)	25 (10%)	7	10
1	B	256/259 (99%)	233 (91%)	23 (9%)	9	13
All	All	512/518 (99%)	464 (91%)	48 (9%)	8	12

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	MET
1	A	16	LYS
1	A	17	LYS
1	A	28	LEU
1	A	31	GLU
1	A	54	TYR
1	A	57	ILE
1	A	58	GLN
1	A	97	LYS
1	A	98	ASN
1	A	118	THR
1	A	140	VAL
1	A	144	ILE
1	A	164	LEU
1	A	170	LEU
1	A	189	LEU
1	A	193	GLU
1	A	199	VAL
1	A	201	GLN
1	A	230	ILE
1	A	232	VAL
1	A	236	GLU
1	A	248	TYR
1	A	275	LYS
1	A	276	LEU
1	B	3	MET
1	B	17	LYS

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Mol	Chain	Res	Type
1	B	53	VAL
1	B	54	TYR
1	B	57	ILE
1	B	60	LEU
1	B	85	ARG
1	B	88	MET
1	B	97	LYS
1	B	98	ASN
1	B	118	THR
1	B	133	VAL
1	B	164	LEU
1	B	170	LEU
1	B	180	GLU
1	B	189	LEU
1	B	232	VAL
1	B	235	ILE
1	B	248	TYR
1	B	257	GLN
1	B	261	THR
1	B	272	LYS
1	B	276	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	68	GLN
1	A	92	HIS
1	A	96	HIS
1	A	102	GLN
1	A	152	GLN
1	A	168	ASN
1	A	187	GLN
1	A	191	GLN
1	A	201	GLN
1	A	207	HIS
1	A	238	GLN
1	A	257	GLN
1	B	7	ASN
1	B	61	ASN
1	B	81	GLN
1	B	115	GLN

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Mol	Chain	Res	Type
1	B	148	HIS
1	B	152	GLN
1	B	191	GLN
1	B	207	HIS
1	B	238	GLN
1	B	259	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 10 ligands modelled in this entry, 6 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	B	1003	3	25,28,28	2.98	3 (12%)	30,37,37	1.74	10 (33%)
2	GGs	A	1002	3	25,28,28	2.93	3 (12%)	30,37,37	1.69	7 (23%)
2	GGs	B	1004	3	25,28,28	2.91	2 (8%)	30,37,37	1.73	5 (16%)
2	GGs	A	1001	3	25,28,28	3.00	3 (12%)	30,37,37	2.02	12 (40%)

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	B	1003	3	-	17/25/31/31	-
2	GGS	A	1002	3	-	17/25/31/31	-
2	GGS	B	1004	3	-	13/25/31/31	-
2	GGS	A	1001	3	-	18/25/31/31	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	GGS	C1-S1	-14.32	1.66	1.84
2	B	1003	GGS	C1-S1	-14.21	1.66	1.84
2	A	1002	GGS	C1-S1	-13.95	1.67	1.84
2	B	1004	GGS	C1-S1	-13.86	1.67	1.84
2	B	1003	GGS	PA-O1A	2.29	1.51	1.46
2	B	1004	GGS	PA-O2A	-2.21	1.51	1.56
2	A	1002	GGS	PA-O2A	-2.20	1.51	1.56
2	A	1001	GGS	PA-O1A	2.13	1.51	1.46
2	A	1001	GGS	PA-O2A	-2.12	1.51	1.56
2	A	1002	GGS	PA-O1A	2.11	1.51	1.46
2	B	1003	GGS	PA-O2A	-2.02	1.51	1.56

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1002	GGS	C10-C8-C9	5.25	124.34	115.23
2	A	1001	GGS	C11-C12-C13	-4.44	117.46	127.62
2	B	1004	GGS	C10-C8-C9	4.39	122.85	115.23
2	A	1001	GGS	C10-C8-C9	3.93	122.05	115.23
2	A	1001	GGS	C14-C13-C15	3.52	121.34	115.23
2	B	1003	GGS	C10-C8-C9	3.39	121.11	115.23
2	B	1003	GGS	C4-C3-C5	3.11	120.63	115.23
2	A	1002	GGS	C2-C1-S1	3.06	126.82	111.87
2	B	1004	GGS	C15-C16-C17	2.88	126.26	112.02
2	B	1004	GGS	C6-C7-C8	-2.82	121.17	127.62
2	A	1002	GGS	C10-C8-C7	-2.82	116.39	123.63
2	B	1003	GGS	C11-C12-C13	-2.81	121.18	127.62
2	A	1001	GGS	C6-C5-C3	-2.60	104.59	113.19
2	A	1001	GGS	C6-C7-C8	-2.56	121.75	127.62
2	A	1002	GGS	C14-C13-C15	2.55	119.66	115.23
2	B	1003	GGS	C2-C1-S1	2.55	124.32	111.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1004	GGs	C16-C17-C18	-2.40	119.66	127.64
2	B	1003	GGs	C15-C16-C17	2.40	123.87	112.02
2	A	1001	GGs	O3A-PB-O1B	-2.36	98.65	111.04
2	A	1002	GGs	C14-C13-C12	-2.35	117.58	123.63
2	A	1002	GGs	O3B-PB-O2B	2.35	116.62	107.80
2	A	1001	GGs	O2A-PA-O1A	2.31	115.36	109.86
2	A	1001	GGs	C15-C16-C17	2.29	123.33	112.02
2	A	1001	GGs	C14-C13-C12	-2.28	117.78	123.63
2	B	1003	GGs	C5-C6-C7	2.26	123.20	112.02
2	B	1003	GGs	C4-C3-C2	-2.21	117.96	123.63
2	A	1001	GGs	O3B-PB-O2B	2.19	116.02	107.80
2	B	1003	GGs	C6-C5-C3	-2.18	105.96	113.19
2	B	1003	GGs	O2A-PA-O1A	2.14	114.97	109.86
2	B	1003	GGs	C14-C13-C15	2.10	118.87	115.23
2	B	1004	GGs	C2-C1-S1	2.08	122.00	111.87
2	A	1001	GGs	C5-C6-C7	2.07	122.24	112.02
2	A	1002	GGs	O2A-PA-O1A	2.06	114.77	109.86
2	A	1001	GGs	C2-C1-S1	2.05	121.87	111.87

There are no chirality outliers.

All (65) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1001	GGs	C1-C2-C3-C4
2	A	1001	GGs	C1-C2-C3-C5
2	A	1001	GGs	C5-C6-C7-C8
2	A	1001	GGs	C6-C7-C8-C9
2	A	1001	GGs	C6-C7-C8-C10
2	A	1001	GGs	PA-O3A-PB-O2B
2	A	1001	GGs	C15-C16-C17-C18
2	A	1001	GGs	C16-C17-C18-C19
2	A	1002	GGs	C6-C7-C8-C9
2	A	1002	GGs	C6-C7-C8-C10
2	A	1002	GGs	C16-C17-C18-C19
2	B	1003	GGs	C1-C2-C3-C4
2	B	1003	GGs	C1-C2-C3-C5
2	B	1003	GGs	C5-C6-C7-C8
2	B	1003	GGs	PA-O3A-PB-O2B
2	B	1003	GGs	C15-C16-C17-C18
2	B	1003	GGs	C16-C17-C18-C20
2	B	1004	GGs	C1-C2-C3-C5
2	B	1004	GGs	C11-C12-C13-C14

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Mol	Chain	Res	Type	Atoms
2	A	1001	GGG	C16-C17-C18-C20
2	A	1002	GGG	C16-C17-C18-C20
2	B	1003	GGG	C16-C17-C18-C19
2	B	1004	GGG	C16-C17-C18-C19
2	B	1004	GGG	C16-C17-C18-C20
2	A	1002	GGG	C10-C8-C9-C11
2	A	1002	GGG	C14-C13-C15-C16
2	B	1003	GGG	C4-C3-C5-C6
2	A	1002	GGG	C12-C13-C15-C16
2	B	1003	GGG	C2-C3-C5-C6
2	A	1002	GGG	C11-C12-C13-C14
2	B	1003	GGG	C6-C7-C8-C10
2	A	1002	GGG	C11-C12-C13-C15
2	B	1003	GGG	C6-C7-C8-C9
2	B	1004	GGG	C11-C12-C13-C15
2	A	1002	GGG	C7-C8-C9-C11
2	A	1001	GGG	C3-C5-C6-C7
2	A	1001	GGG	C12-C11-C9-C8
2	A	1002	GGG	C12-C11-C9-C8
2	A	1002	GGG	C13-C15-C16-C17
2	B	1003	GGG	C3-C5-C6-C7
2	B	1004	GGG	C10-C8-C9-C11
2	B	1004	GGG	C7-C8-C9-C11
2	B	1004	GGG	C12-C11-C9-C8
2	A	1001	GGG	C4-C3-C5-C6
2	A	1002	GGG	C3-C5-C6-C7
2	B	1003	GGG	PA-O3A-PB-O1B
2	B	1004	GGG	C1-C2-C3-C4
2	A	1001	GGG	C2-C3-C5-C6
2	B	1003	GGG	C10-C8-C9-C11
2	A	1002	GGG	PA-O3A-PB-O2B
2	A	1001	GGG	C10-C8-C9-C11
2	B	1003	GGG	C7-C8-C9-C11
2	A	1001	GGG	C7-C8-C9-C11
2	A	1001	GGG	PB-O3A-PA-O2A
2	B	1003	GGG	PB-O3A-PA-O2A
2	B	1004	GGG	PB-O3A-PA-O1A
2	A	1001	GGG	PA-O3A-PB-O1B
2	A	1001	GGG	C9-C11-C12-C13
2	A	1002	GGG	C5-C6-C7-C8
2	A	1002	GGG	C15-C16-C17-C18
2	B	1003	GGG	C9-C11-C12-C13

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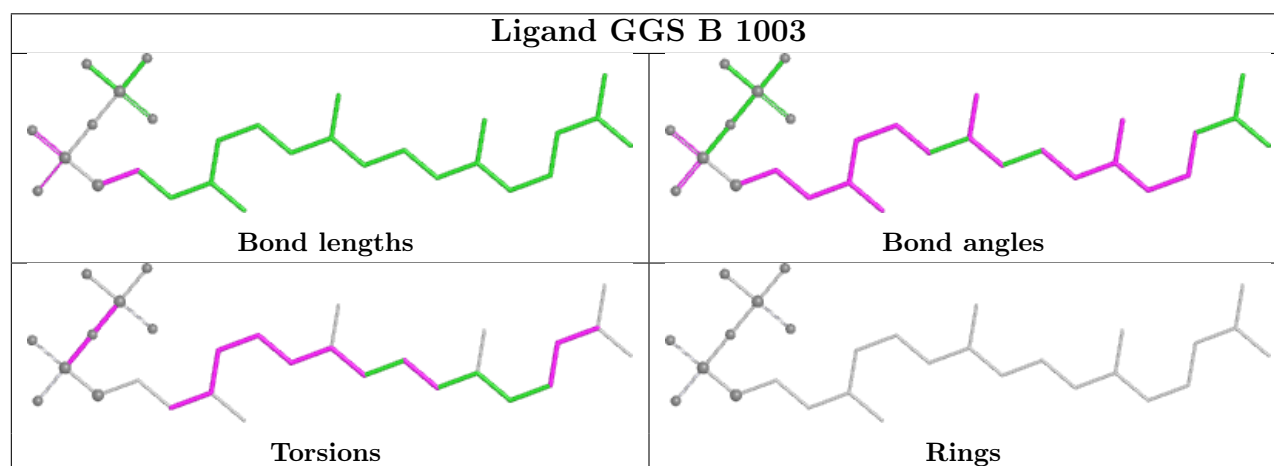
Mol	Chain	Res	Type	Atoms
2	A	1002	GGs	C9-C11-C12-C13
2	B	1004	GGs	C5-C6-C7-C8
2	B	1004	GGs	C9-C11-C12-C13
2	B	1004	GGs	C15-C16-C17-C18

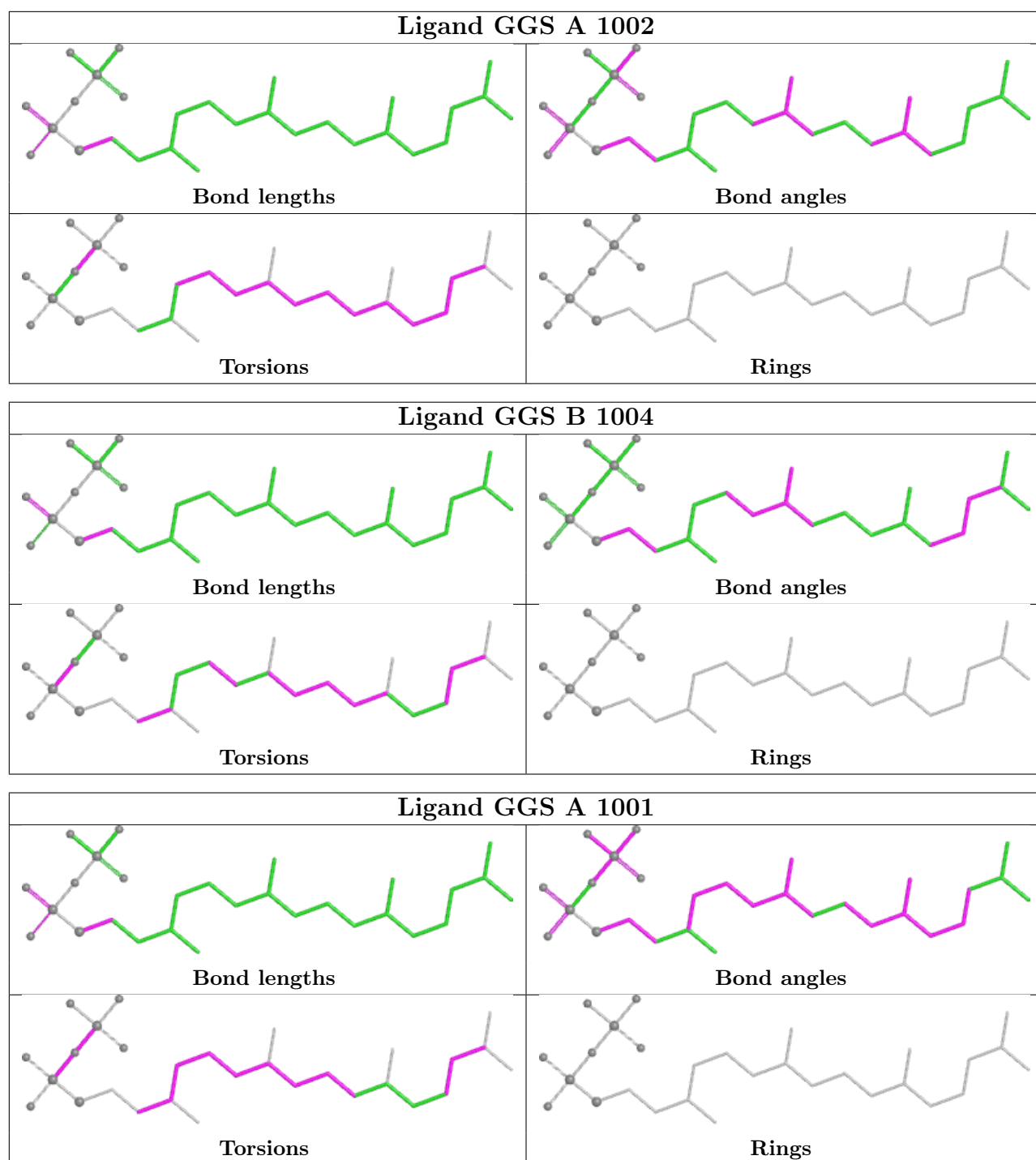
There are no ring outliers.

4 monomers are involved in 58 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1003	GGs	20	0
2	A	1002	GGs	20	0
2	B	1004	GGs	18	0
2	A	1001	GGs	21	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, 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	284/293 (96%)	0.37	8 (2%) 55 54	15, 29, 42, 56	0
1	B	284/293 (96%)	0.47	15 (5%) 32 31	19, 29, 47, 61	0
All	All	568/586 (96%)	0.42	23 (4%) 42 42	15, 29, 44, 61	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	17	LYS	3.4
1	B	56	ASP	3.3
1	A	202	ASN	3.0
1	A	17	LYS	2.9
1	B	200	TYR	2.7
1	A	163	SER	2.6
1	A	43	VAL	2.5
1	B	73	TYR	2.5
1	A	282	SER	2.4
1	A	200	TYR	2.4
1	B	51	ILE	2.3
1	B	198	GLU	2.2
1	B	96	HIS	2.2
1	B	202	ASN	2.2
1	A	116	HIS	2.2
1	B	259	ASN	2.1
1	B	203	GLY	2.1
1	B	50	SER	2.1
1	B	54	TYR	2.1
1	B	57	ILE	2.1
1	B	115	GLN	2.1
1	A	177	PHE	2.1
1	B	267	PHE	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

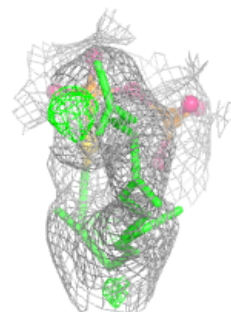
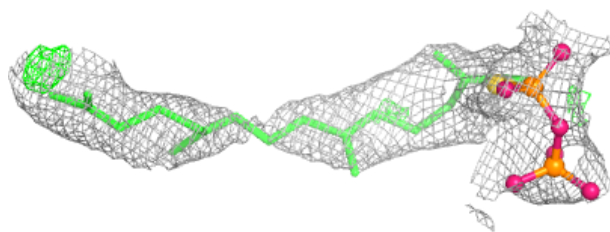
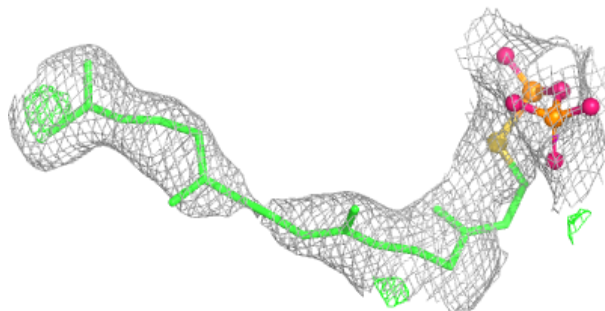
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MG	B	1010	1/1	0.78	0.13	67,67,67,67	0
3	MG	B	1009	1/1	0.84	0.14	59,59,59,59	0
3	MG	A	1006	1/1	0.84	0.07	55,55,55,55	0
2	GGs	B	1004	29/29	0.86	0.14	60,69,80,82	0
3	MG	A	1007	1/1	0.86	0.17	65,65,65,65	0
2	GGs	A	1002	29/29	0.88	0.14	53,60,68,70	0
2	GGs	A	1001	29/29	0.88	0.16	48,70,75,76	0
2	GGs	B	1003	29/29	0.90	0.15	52,70,77,77	0
3	MG	B	1008	1/1	0.92	0.14	44,44,44,44	0
3	MG	A	1005	1/1	0.97	0.13	42,42,42,42	0

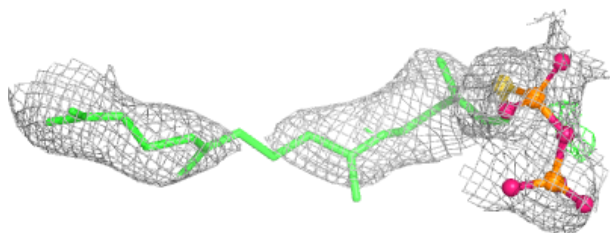
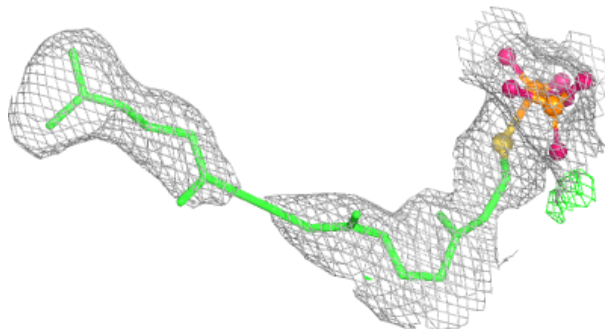
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around GGS B 1004:

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

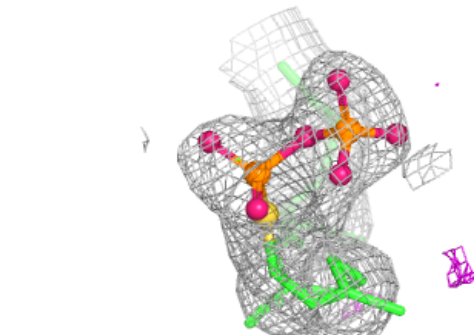
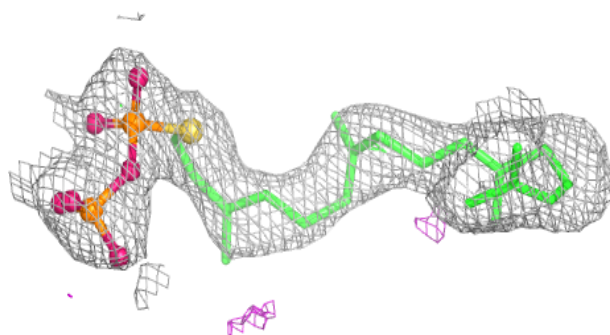
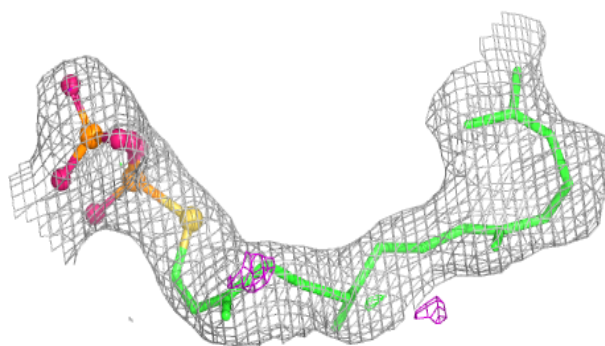
**Electron density around GGS A 1002:**

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

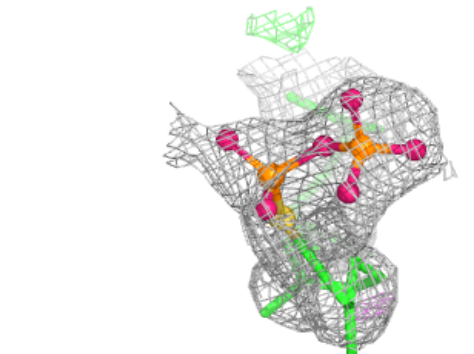
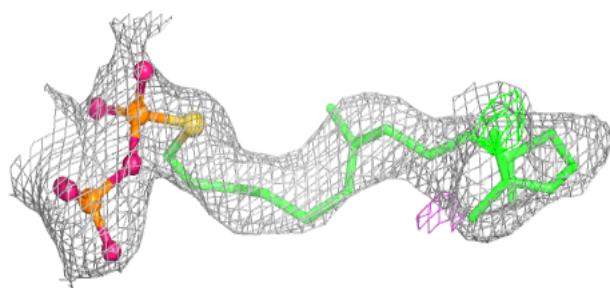
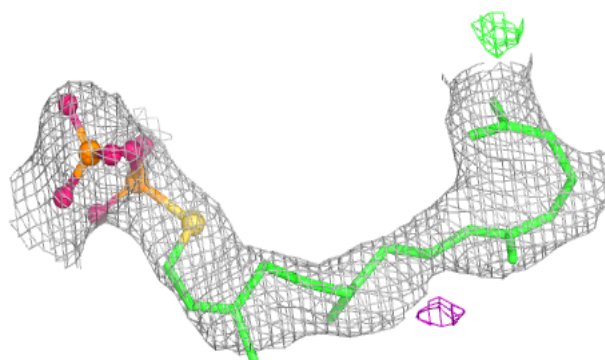


Electron density around GGS A 1001:

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

**Electron density around GGS B 1003:**

$2mF_o-DF_c$ (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.