



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 12:40 PM UTC

PDB ID : 9ZAV / pdb_00009zav
Title : Crystal structure of Formyl-coenzyme A transferase from Brucella melitensis in complex with succinate
Authors : Seattle Structural Genomics Center for Infectious Disease; Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2025-11-19
Resolution : 2.88 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

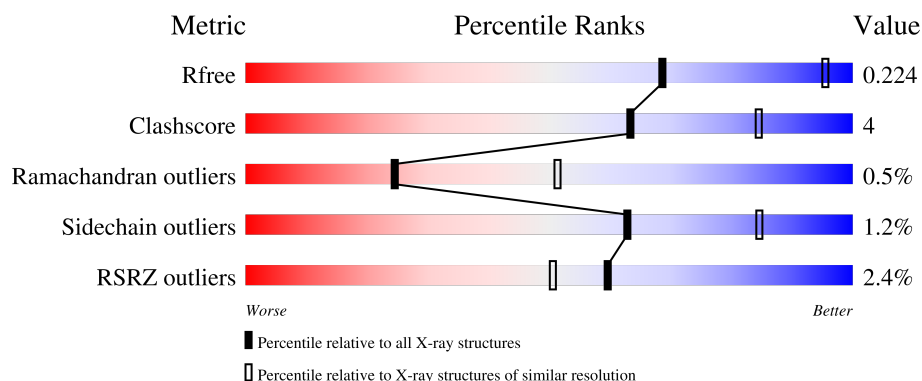
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.88 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	3557 (2.90-2.86)
Clashscore	190562	3801 (2.90-2.86)
Ramachandran outliers	187476	3699 (2.90-2.86)
Sidechain outliers	187428	3702 (2.90-2.86)
RSRZ outliers	180081	3558 (2.90-2.86)

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	409	
1	B	409	
1	C	409	
1	D	409	

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Mol	Chain	Length	Quality of chain
1	E	409	 A horizontal bar chart showing the quality of chain E. The bar is divided into four segments: a small red segment at the beginning labeled '6%', a large green segment labeled '86%', a small yellow segment labeled '11%', and a very small grey segment at the end. A small black dot is located at the far right end of the bar.

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 15092 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 Formyl-coenzyme a transferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	403	Total	C	N	O	S	0	0	0
			3044	1921	535	573	15			
1	B	402	Total	C	N	O	S	0	0	0
			3030	1912	531	572	15			
1	C	402	Total	C	N	O	S	0	0	0
			3026	1910	531	570	15			
1	D	402	Total	C	N	O	S	0	0	0
			2999	1893	520	571	15			
1	E	397	Total	C	N	O	S	0	0	0
			2969	1877	515	563	14			

There are 40 discrepancies between the modelled and reference sequences:

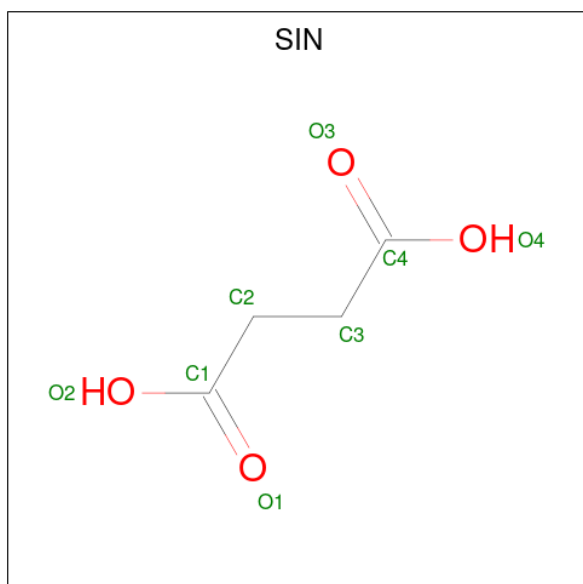
Chain	Residue	Modelled	Actual	Comment	Reference
A	42	MET	-	expression tag	UNP Q8YDF2
A	43	ALA	-	expression tag	UNP Q8YDF2
A	44	HIS	-	expression tag	UNP Q8YDF2
A	45	HIS	-	expression tag	UNP Q8YDF2
A	46	HIS	-	expression tag	UNP Q8YDF2
A	47	HIS	-	expression tag	UNP Q8YDF2
A	48	HIS	-	expression tag	UNP Q8YDF2
A	49	HIS	-	expression tag	UNP Q8YDF2
B	42	MET	-	expression tag	UNP Q8YDF2
B	43	ALA	-	expression tag	UNP Q8YDF2
B	44	HIS	-	expression tag	UNP Q8YDF2
B	45	HIS	-	expression tag	UNP Q8YDF2
B	46	HIS	-	expression tag	UNP Q8YDF2
B	47	HIS	-	expression tag	UNP Q8YDF2
B	48	HIS	-	expression tag	UNP Q8YDF2
B	49	HIS	-	expression tag	UNP Q8YDF2
C	42	MET	-	expression tag	UNP Q8YDF2
C	43	ALA	-	expression tag	UNP Q8YDF2
C	44	HIS	-	expression tag	UNP Q8YDF2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	45	HIS	-	expression tag	UNP Q8YDF2
C	46	HIS	-	expression tag	UNP Q8YDF2
C	47	HIS	-	expression tag	UNP Q8YDF2
C	48	HIS	-	expression tag	UNP Q8YDF2
C	49	HIS	-	expression tag	UNP Q8YDF2
D	42	MET	-	expression tag	UNP Q8YDF2
D	43	ALA	-	expression tag	UNP Q8YDF2
D	44	HIS	-	expression tag	UNP Q8YDF2
D	45	HIS	-	expression tag	UNP Q8YDF2
D	46	HIS	-	expression tag	UNP Q8YDF2
D	47	HIS	-	expression tag	UNP Q8YDF2
D	48	HIS	-	expression tag	UNP Q8YDF2
D	49	HIS	-	expression tag	UNP Q8YDF2
E	42	MET	-	expression tag	UNP Q8YDF2
E	43	ALA	-	expression tag	UNP Q8YDF2
E	44	HIS	-	expression tag	UNP Q8YDF2
E	45	HIS	-	expression tag	UNP Q8YDF2
E	46	HIS	-	expression tag	UNP Q8YDF2
E	47	HIS	-	expression tag	UNP Q8YDF2
E	48	HIS	-	expression tag	UNP Q8YDF2
E	49	HIS	-	expression tag	UNP Q8YDF2

- Molecule 2 is SUCCINIC ACID (CCD ID: SIN) (formula: $C_4H_6O_4$) (labeled as "Ligand of Interest" by depositor).



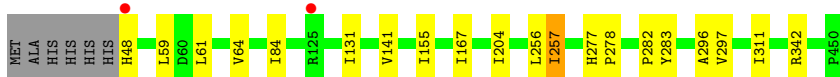
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 8	C 4	O 4	0	0
2	B	1	Total 8	C 4	O 4	0	0
2	C	1	Total 8	C 4	O 4	0	0

3 Residue-property plots

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.

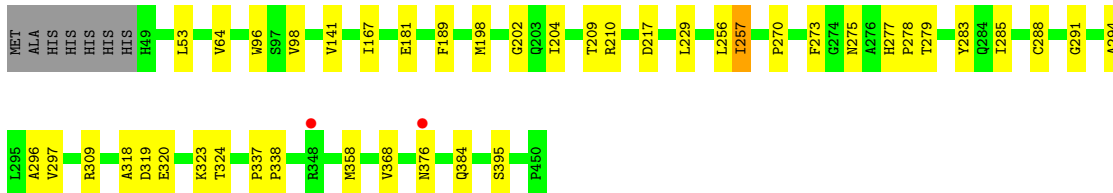
- Molecule 1: Formyl-coenzyme a transferase

Chain A: 



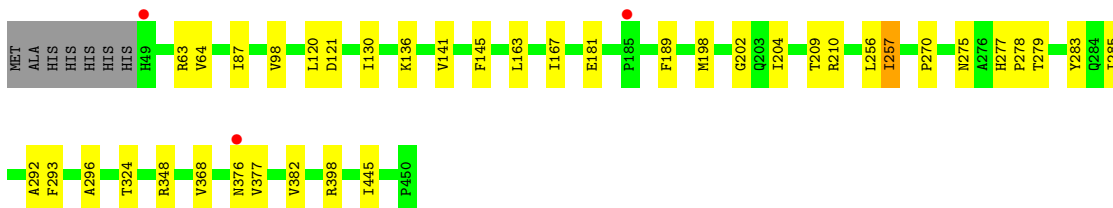
- Molecule 1: Formyl-coenzyme a transferase

Chain B: 



- Molecule 1: Formyl-coenzyme a transferase

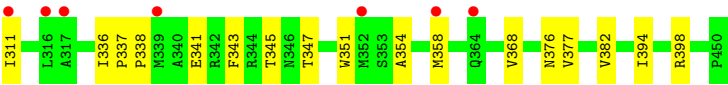
Chain C: 



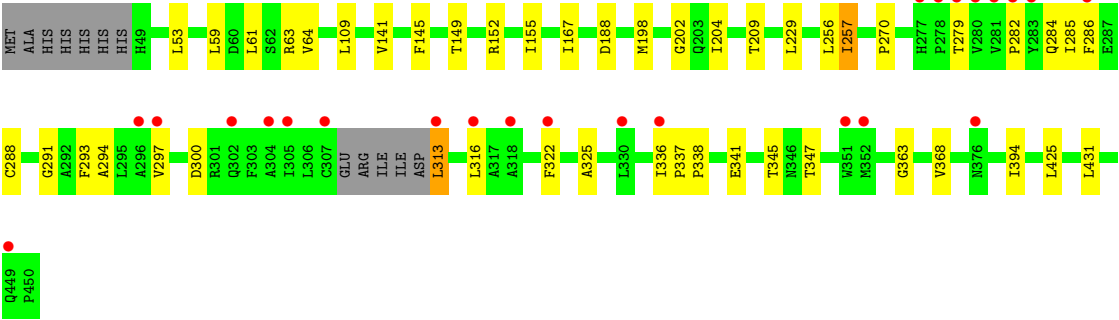
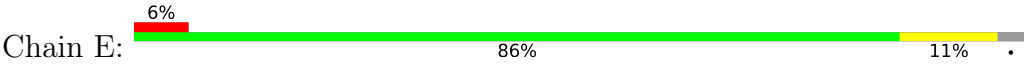
- Molecule 1: Formyl-coenzyme a transferase

Chain D: 





● Molecule 1: Formyl-coenzyme a transferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	189.59Å 189.59Å 199.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.19 – 2.88 48.19 – 2.88	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.19-2.88) 100.0 (48.19-2.88)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 2.86Å)	Xtriage
Refinement program	PHENIX (2.0_5819: ???)	Depositor
R, R_{free}	0.192 , 0.225 0.194 , 0.224	Depositor DCC
R_{free} test set	4273 reflections (5.18%)	wwPDB-VP
Wilson B-factor (Å ²)	79.7	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 67.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k 0.000 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15092	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.12% 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, 0.333 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: SIN

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	0.17	0/3103	0.38	0/4211
1	B	0.17	0/3088	0.35	0/4192
1	C	0.16	0/3084	0.35	0/4187
1	D	0.16	0/3057	0.34	0/4155
1	E	0.15	0/3026	0.33	0/4110
All	All	0.17	0/15358	0.35	0/20855

There are no bond length outliers.

There are no bond angle outliers.

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	3044	0	3055	13	0
1	B	3030	0	3037	40	0
1	C	3026	0	3033	34	0
1	D	2999	0	2971	28	0
1	E	2969	0	2962	31	0
2	A	8	0	4	0	0
2	B	8	0	4	1	0
2	C	8	0	4	0	0
All	All	15092	0	15070	119	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 (119) 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:D:285:ILE:HD11	1:E:198:MET:HE1	1.34	1.08
1:D:198:MET:HE1	1:E:285:ILE:HD11	1.49	0.94
1:A:141:VAL:HG23	1:A:167:ILE:HD11	1.64	0.78
1:B:319:ASP:O	1:B:323:LYS:HG3	1.84	0.75
1:E:284:GLN:NE2	1:E:286:PHE:CZ	2.55	0.74
1:E:141:VAL:HG23	1:E:167:ILE:HD11	1.72	0.72
1:E:282:PRO:HD2	1:E:297:VAL:HG21	1.71	0.71
1:B:285:ILE:HD11	1:C:198:MET:HE1	1.76	0.67
1:B:318:ALA:O	1:B:323:LYS:HE2	1.96	0.66
1:B:198:MET:HE1	1:C:285:ILE:HD11	1.83	0.61
1:D:198:MET:CE	1:E:285:ILE:HD11	2.29	0.60
1:C:141:VAL:HG23	1:C:167:ILE:HD11	1.83	0.58
1:B:320:GLU:N	1:B:323:LYS:HE3	2.19	0.58
1:C:141:VAL:CG2	1:C:167:ILE:HD11	2.33	0.58
1:D:310:ILE:HG22	1:D:351:TRP:CE3	2.40	0.57
1:B:279:THR:HG23	1:C:98:VAL:O	2.05	0.56
1:B:204:ILE:H	1:B:204:ILE:HD12	1.69	0.56
1:E:336:ILE:HG13	1:E:337:PRO:HD3	1.86	0.56
1:B:320:GLU:HA	1:B:323:LYS:HG3	1.86	0.56
1:D:377:VAL:HG13	1:D:382:VAL:HB	1.86	0.56
1:E:313:LEU:HD23	1:E:316:LEU:HD11	1.87	0.56
1:D:204:ILE:HD12	1:D:204:ILE:H	1.71	0.56
1:A:141:VAL:CG2	1:A:167:ILE:HD11	2.36	0.54
1:C:87:ILE:HD11	1:C:121:ASP:HB2	1.89	0.54
1:C:120:LEU:CD1	1:C:445:ILE:HG22	2.38	0.54
1:E:336:ILE:HD12	1:E:337:PRO:N	2.22	0.54
1:D:341:GLU:O	1:D:345:THR:HG23	2.08	0.54
1:D:336:ILE:HG13	1:D:337:PRO:HD3	1.90	0.53
1:B:318:ALA:O	1:B:323:LYS:CE	2.56	0.53
1:E:61:LEU:HD13	1:E:155:ILE:HD12	1.90	0.53
1:C:120:LEU:HD12	1:C:445:ILE:HG22	1.89	0.53
1:D:270:PRO:HG2	1:E:368:VAL:HB	1.90	0.53
1:A:61:LEU:HD13	1:A:155:ILE:HD12	1.90	0.53
1:D:59:LEU:HB3	1:D:141:VAL:HG22	1.91	0.52
1:B:294:ALA:HB1	1:C:189:PHE:CE2	2.44	0.52
1:E:204:ILE:HD12	1:E:204:ILE:H	1.75	0.52
1:B:181:GLU:HB2	1:C:376:ASN:OD1	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:ILE:HD12	1:A:204:ILE:H	1.76	0.51
1:B:309:ARG:HB3	1:B:358:MET:HE3	1.92	0.51
1:E:279:THR:HG22	1:E:279:THR:O	2.11	0.51
1:B:210:ARG:HG2	1:C:275:ASN:HA	1.93	0.51
1:D:209:THR:HG21	1:E:202:GLY:HA2	1.93	0.50
1:C:204:ILE:H	1:C:204:ILE:HD12	1.75	0.50
1:D:279:THR:O	1:D:279:THR:HG22	2.11	0.50
1:A:282:PRO:HD2	1:A:297:VAL:HG21	1.93	0.50
1:B:368:VAL:HB	1:C:270:PRO:HG2	1.92	0.49
1:D:368:VAL:HB	1:E:270:PRO:HG2	1.94	0.49
1:E:341:GLU:O	1:E:345:THR:HG23	2.11	0.49
1:D:284:GLN:HE22	1:D:336:ILE:HG21	1.77	0.48
1:B:320:GLU:HA	1:B:323:LYS:CD	2.44	0.48
1:A:311:ILE:HB	1:A:342:ARG:HG2	1.95	0.48
1:B:53:LEU:HD11	1:B:229:LEU:HD21	1.95	0.48
1:D:190:VAL:HG13	1:E:363:GLY:HA3	1.95	0.48
1:E:337:PRO:HB2	1:E:338:PRO:HD3	1.96	0.48
1:C:277:HIS:CG	1:C:278:PRO:HD2	2.49	0.47
1:B:283:TYR:CE1	1:B:296:ALA:HB1	2.49	0.47
1:C:87:ILE:CD1	1:C:121:ASP:HB2	2.43	0.47
1:B:270:PRO:HG2	1:C:368:VAL:HB	1.97	0.47
1:E:288:CYS:HB2	1:E:291:GLY:O	2.15	0.47
1:D:202:GLY:HA2	1:E:209:THR:HG21	1.96	0.46
1:B:320:GLU:CA	1:B:323:LYS:HE3	2.46	0.46
1:D:354:ALA:O	1:D:358:MET:HG2	2.15	0.46
1:B:285:ILE:HG13	1:C:189:PHE:HZ	1.81	0.46
1:B:275:ASN:HA	1:C:210:ARG:HG2	1.96	0.46
1:C:136:LYS:HE2	1:C:163:LEU:HD22	1.98	0.46
1:E:313:LEU:HD23	1:E:316:LEU:CD1	2.45	0.46
1:C:120:LEU:HD11	1:C:130:ILE:HD12	1.98	0.45
1:A:283:TYR:CE1	1:A:296:ALA:HB1	2.51	0.45
1:C:283:TYR:CE1	1:C:296:ALA:HB1	2.51	0.45
1:E:53:LEU:HD11	1:E:229:LEU:HD21	1.97	0.45
1:B:98:VAL:O	1:C:279:THR:HG23	2.16	0.45
1:D:310:ILE:HG22	1:D:351:TRP:CZ3	2.51	0.45
1:B:285:ILE:CD1	1:C:198:MET:HE1	2.46	0.45
1:D:336:ILE:HD12	1:D:337:PRO:N	2.32	0.45
1:E:59:LEU:HB3	1:E:141:VAL:HG22	1.97	0.45
1:E:300:ASP:OD2	1:E:325:ALA:HB2	2.17	0.45
1:B:273:PHE:CE1	1:B:278:PRO:HB3	2.52	0.45
1:C:293:PHE:HB3	1:C:348:ARG:HG3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:376:ASN:OD1	1:C:181:GLU:HB2	2.17	0.45
1:A:84:ILE:HG21	1:A:131:ILE:HD11	1.98	0.44
1:B:202:GLY:HA2	1:C:209:THR:HG21	1.98	0.44
1:C:377:VAL:HG13	1:C:382:VAL:HB	1.99	0.44
1:B:277:HIS:CG	1:B:278:PRO:HD2	2.53	0.44
1:C:285:ILE:HG23	1:C:292:ALA:HB1	2.00	0.44
1:B:297:VAL:HG23	1:B:297:VAL:O	2.18	0.43
1:D:189:PHE:CE2	1:E:294:ALA:HB1	2.53	0.43
1:E:282:PRO:O	1:E:297:VAL:HG22	2.18	0.43
1:B:256:LEU:O	1:B:257:ILE:C	2.61	0.43
1:D:109:LEU:HD22	1:D:394:ILE:HD13	2.00	0.43
1:D:336:ILE:N	1:D:337:PRO:CD	2.81	0.43
1:D:337:PRO:HB2	1:D:338:PRO:HD3	1.99	0.43
1:A:256:LEU:O	1:A:257:ILE:C	2.61	0.43
1:B:288:CYS:HB2	1:B:291:GLY:O	2.18	0.43
1:B:320:GLU:HA	1:B:323:LYS:CG	2.48	0.43
1:A:297:VAL:HG23	1:A:297:VAL:O	2.16	0.43
1:B:320:GLU:HA	1:B:323:LYS:HE3	2.01	0.43
1:B:384:GLN:O	1:B:395:SER:HA	2.19	0.43
1:D:282:PRO:HD2	1:D:297:VAL:HG21	1.99	0.43
1:B:337:PRO:HB2	1:B:338:PRO:HD3	2.01	0.43
1:D:284:GLN:NE2	1:D:336:ILE:HG21	2.33	0.43
1:B:96:TRP:HB3	1:C:279:THR:CG2	2.49	0.42
1:A:282:PRO:O	1:A:297:VAL:HG22	2.18	0.42
1:E:109:LEU:HD22	1:E:394:ILE:HD13	2.00	0.42
1:B:189:PHE:HZ	1:C:285:ILE:HG13	1.84	0.42
1:B:198:MET:HE1	1:C:285:ILE:CD1	2.48	0.42
1:E:336:ILE:N	1:E:337:PRO:CD	2.83	0.42
1:A:59:LEU:HB3	1:A:141:VAL:HG22	2.01	0.42
1:D:306:LEU:HD12	1:D:306:LEU:C	2.45	0.41
1:B:217:ASP:OD2	2:B:501:SIN:O1	2.39	0.41
1:E:145:PHE:HB3	1:E:149:THR:HB	2.02	0.41
1:E:256:LEU:O	1:E:257:ILE:C	2.64	0.41
1:C:256:LEU:O	1:C:257:ILE:C	2.64	0.41
1:E:425:LEU:HB3	1:E:431:LEU:HD12	2.03	0.41
1:B:209:THR:HG21	1:C:202:GLY:HA2	2.02	0.40
1:C:63:ARG:NH2	1:C:145:PHE:HA	2.37	0.40
1:A:277:HIS:CG	1:A:278:PRO:HD2	2.56	0.40
1:B:141:VAL:HG23	1:B:167:ILE:HD11	2.03	0.40
1:D:167:ILE:O	1:D:241:GLY:HA3	2.22	0.40
1:D:311:ILE:HD11	1:D:343:PHE:CE1	2.57	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	401/409 (98%)	394 (98%)	5 (1%)	2 (0%)	24	51
1	B	400/409 (98%)	396 (99%)	2 (0%)	2 (0%)	24	51
1	C	400/409 (98%)	394 (98%)	4 (1%)	2 (0%)	24	51
1	D	400/409 (98%)	392 (98%)	6 (2%)	2 (0%)	24	51
1	E	393/409 (96%)	387 (98%)	4 (1%)	2 (0%)	24	51
All	All	1994/2045 (98%)	1963 (98%)	21 (1%)	10 (0%)	24	51

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	257	ILE
1	B	257	ILE
1	C	257	ILE
1	D	257	ILE
1	E	257	ILE
1	A	64	VAL
1	B	64	VAL
1	C	64	VAL
1	D	64	VAL
1	E	64	VAL

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	317/322 (98%)	316 (100%)	1 (0%)	86	95
1	B	315/322 (98%)	314 (100%)	1 (0%)	86	95
1	C	314/322 (98%)	312 (99%)	2 (1%)	78	92
1	D	308/322 (96%)	300 (97%)	8 (3%)	40	71
1	E	307/322 (95%)	300 (98%)	7 (2%)	44	74
All	All	1561/1610 (97%)	1542 (99%)	19 (1%)	63	84

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	HIS
1	B	324	THR
1	C	324	THR
1	C	398	ARG
1	D	49	HIS
1	D	288	CYS
1	D	293	PHE
1	D	297	VAL
1	D	306	LEU
1	D	347	THR
1	D	376	ASN
1	D	398	ARG
1	E	63	ARG
1	E	152	ARG
1	E	188	ASP
1	E	293	PHE
1	E	313	LEU
1	E	322	PHE
1	E	347	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	HIS
1	A	234	GLN
1	A	242	GLN
1	A	254	ASN
1	A	263	HIS
1	A	349	GLN
1	A	390	HIS

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Mol	Chain	Res	Type
1	B	177	GLN
1	B	263	HIS
1	B	346	ASN
1	B	349	GLN
1	B	384	GLN
1	B	390	HIS
1	C	268	HIS
1	C	331	ASN
1	C	349	GLN
1	C	390	HIS
1	D	177	GLN
1	D	203	GLN
1	D	263	HIS
1	D	390	HIS
1	E	49	HIS
1	E	242	GLN
1	E	263	HIS
1	E	268	HIS
1	E	390	HIS

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 ⓘ

3 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	SIN	C	501	-	7,7,7	1.35	0	8,8,8	1.63	1 (12%)
2	SIN	B	501	-	7,7,7	1.36	0	8,8,8	1.47	1 (12%)
2	SIN	A	501	-	7,7,7	1.35	0	8,8,8	1.55	3 (37%)

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	SIN	C	501	-	-	3/5/5/5	-
2	SIN	B	501	-	-	4/5/5/5	-
2	SIN	A	501	-	-	4/5/5/5	-

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	SIN	O3-C4-C3	-2.82	114.15	123.09
2	B	501	SIN	O3-C4-C3	-2.42	115.43	123.09
2	A	501	SIN	O3-C4-C3	-2.33	115.71	123.09
2	A	501	SIN	O1-C1-C2	-2.06	116.55	123.09
2	A	501	SIN	C2-C3-C4	2.02	119.02	113.67

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	501	SIN	C1-C2-C3-C4
2	C	501	SIN	C2-C3-C4-O3
2	A	501	SIN	C2-C3-C4-O4
2	A	501	SIN	C2-C3-C4-O3
2	B	501	SIN	C2-C3-C4-O3
2	C	501	SIN	C2-C3-C4-O4
2	B	501	SIN	C2-C3-C4-O4
2	A	501	SIN	O2-C1-C2-C3
2	B	501	SIN	O2-C1-C2-C3

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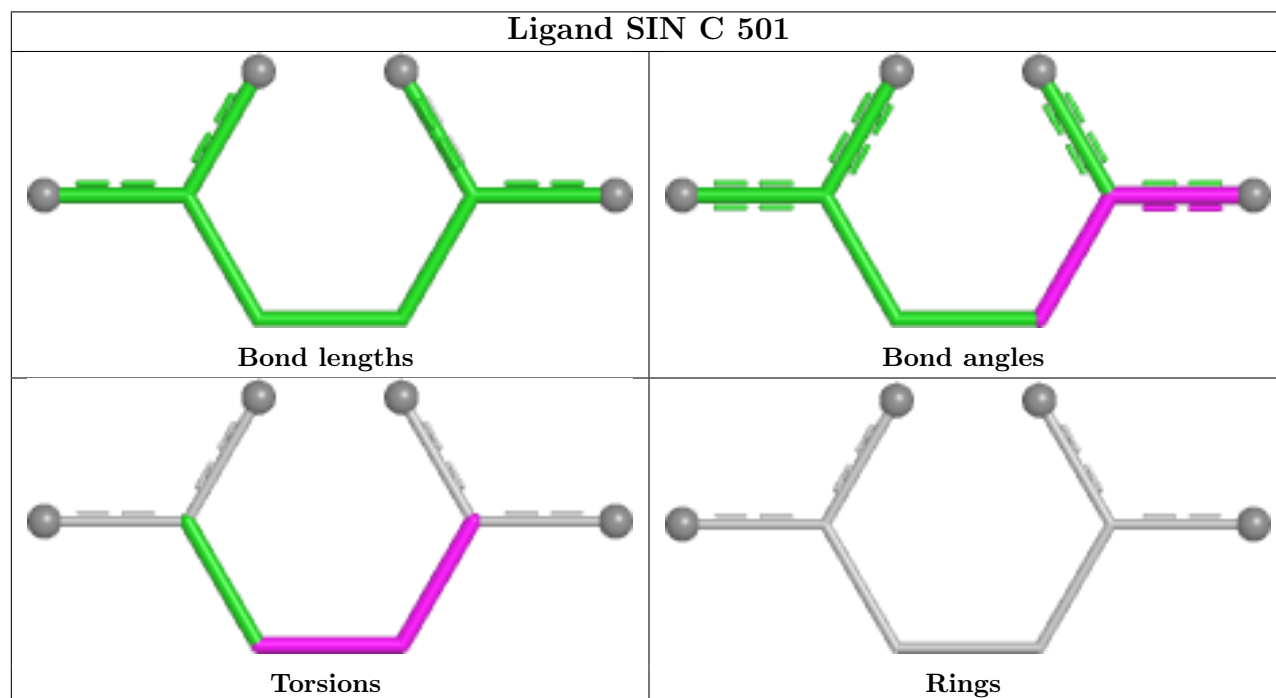
Mol	Chain	Res	Type	Atoms
2	B	501	SIN	O1-C1-C2-C3
2	A	501	SIN	O1-C1-C2-C3

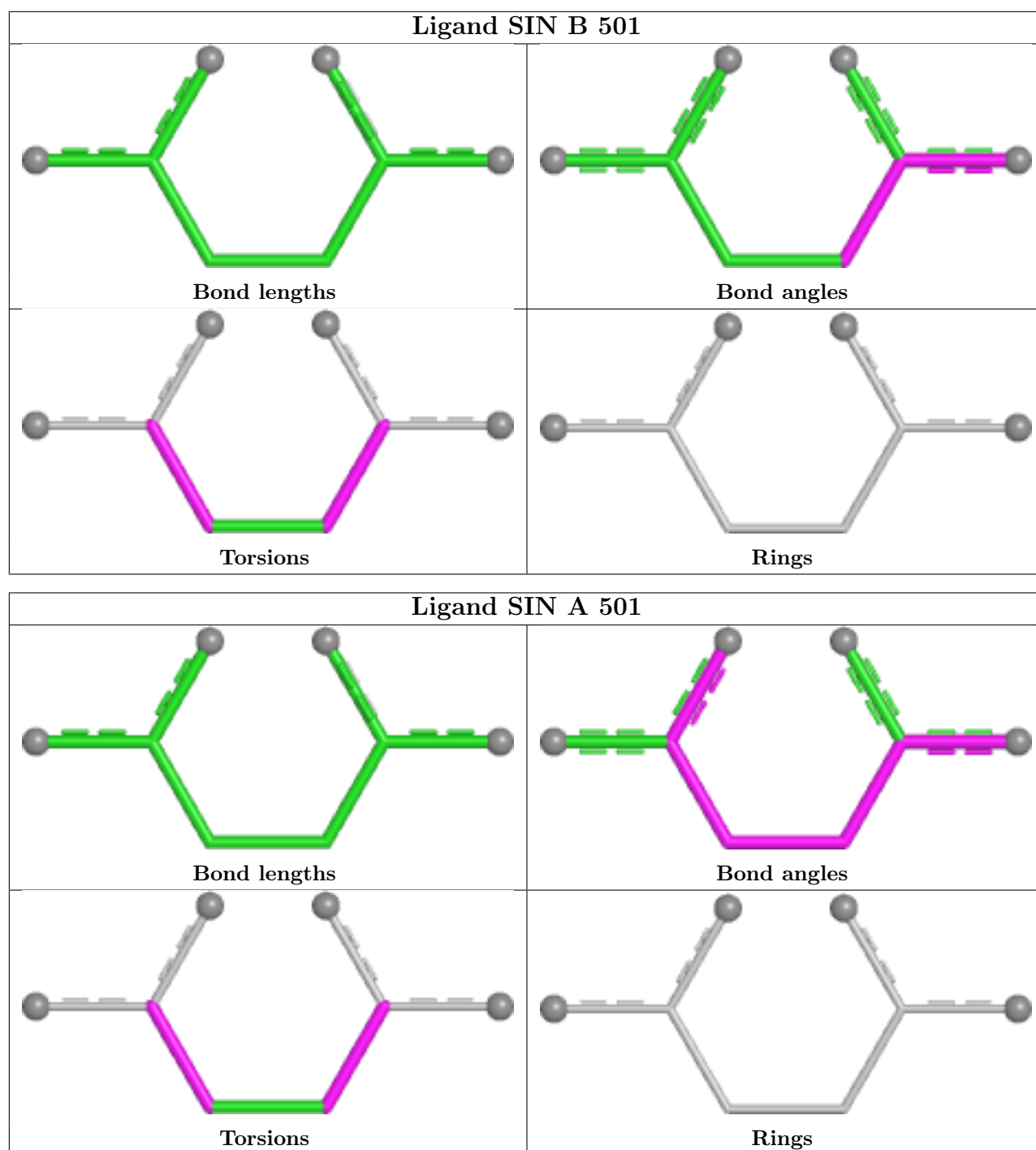
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	SIN	1	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	403/409 (98%)	-0.39	2 (0%) 87 83	44, 64, 92, 146	0
1	B	402/409 (98%)	-0.21	2 (0%) 87 83	47, 74, 110, 136	0
1	C	402/409 (98%)	-0.19	3 (0%) 84 80	48, 75, 112, 132	0
1	D	402/409 (98%)	0.16	17 (4%) 40 32	54, 86, 173, 199	0
1	E	397/409 (97%)	0.17	24 (6%) 27 21	53, 91, 163, 208	0
All	All	2006/2045 (98%)	-0.09	48 (2%) 59 50	44, 77, 153, 208	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	279	THR	4.1
1	D	282	PRO	4.0
1	D	352	MET	3.7
1	E	313	LEU	3.6
1	D	279	THR	3.6
1	E	286	PHE	3.4
1	E	376	ASN	3.2
1	E	304	ALA	3.2
1	A	48	HIS	3.2
1	D	281	VAL	3.1
1	D	305	ILE	3.0
1	E	305	ILE	3.0
1	E	297	VAL	2.9
1	D	317	ALA	2.8
1	C	185	PRO	2.8
1	E	296	ALA	2.8
1	E	281	VAL	2.7
1	E	278	PRO	2.6
1	B	376	ASN	2.6
1	D	364	GLN	2.6

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Mol	Chain	Res	Type	RSRZ
1	E	316	LEU	2.6
1	E	352	MET	2.6
1	E	322	PHE	2.6
1	D	278	PRO	2.6
1	D	268	HIS	2.5
1	E	330	LEU	2.5
1	B	348	ARG	2.5
1	D	316	LEU	2.5
1	A	125	ARG	2.4
1	D	286	PHE	2.4
1	D	277	HIS	2.4
1	E	277	HIS	2.4
1	E	351	TRP	2.4
1	C	49	HIS	2.4
1	D	358	MET	2.4
1	E	302	GLN	2.4
1	D	310	ILE	2.3
1	D	311	ILE	2.3
1	C	376	ASN	2.2
1	E	307	CYS	2.2
1	D	339	MET	2.2
1	E	318	ALA	2.2
1	E	282	PRO	2.2
1	D	296	ALA	2.1
1	E	336	ILE	2.1
1	E	283	TYR	2.1
1	E	449	GLN	2.0
1	E	280	VAL	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 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,

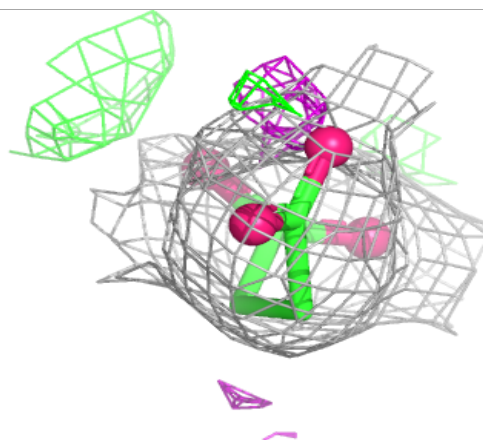
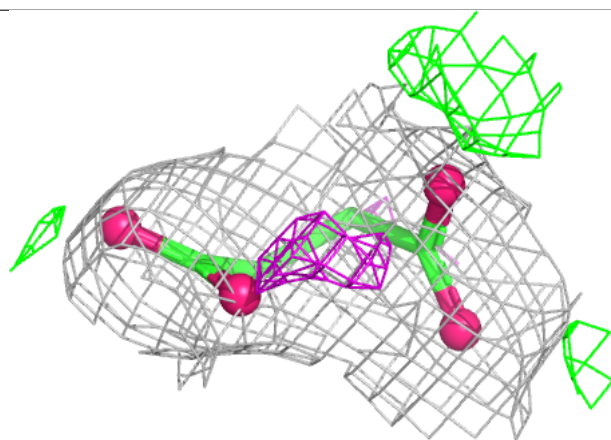
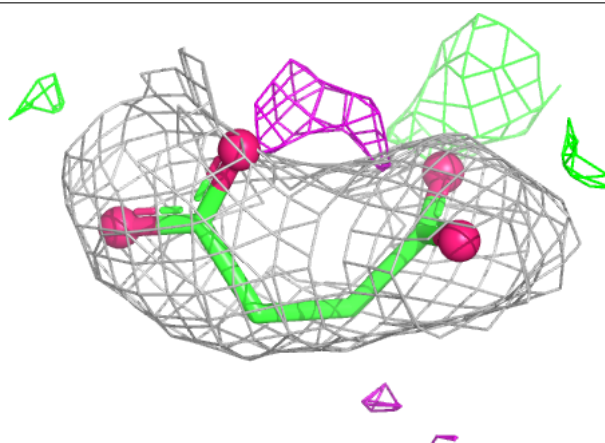
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SIN	A	501	8/8	0.92	0.15	59,70,73,77	0
2	SIN	B	501	8/8	0.92	0.16	67,73,82,82	0
2	SIN	C	501	8/8	0.92	0.16	67,73,78,82	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.

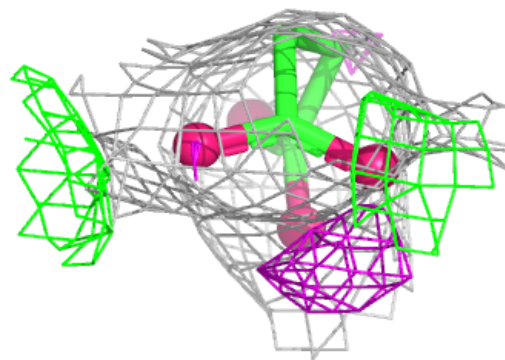
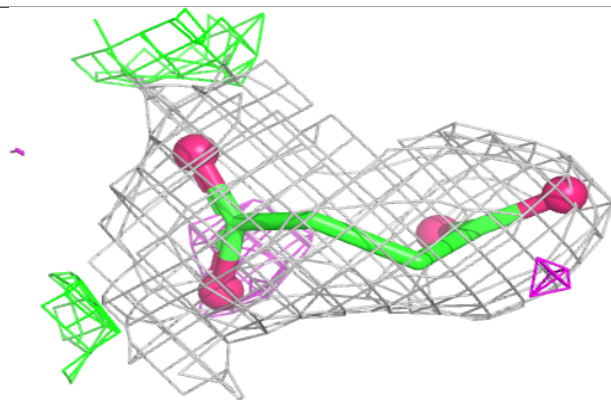
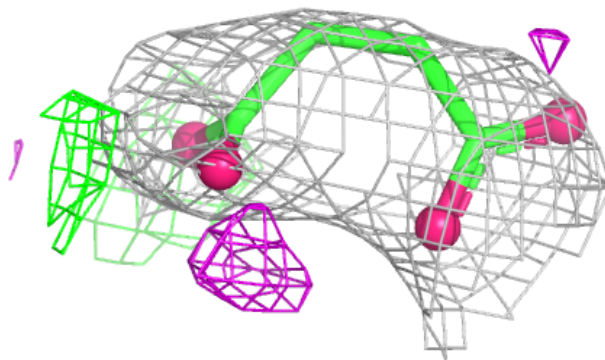
Electron density around SIN A 501:

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



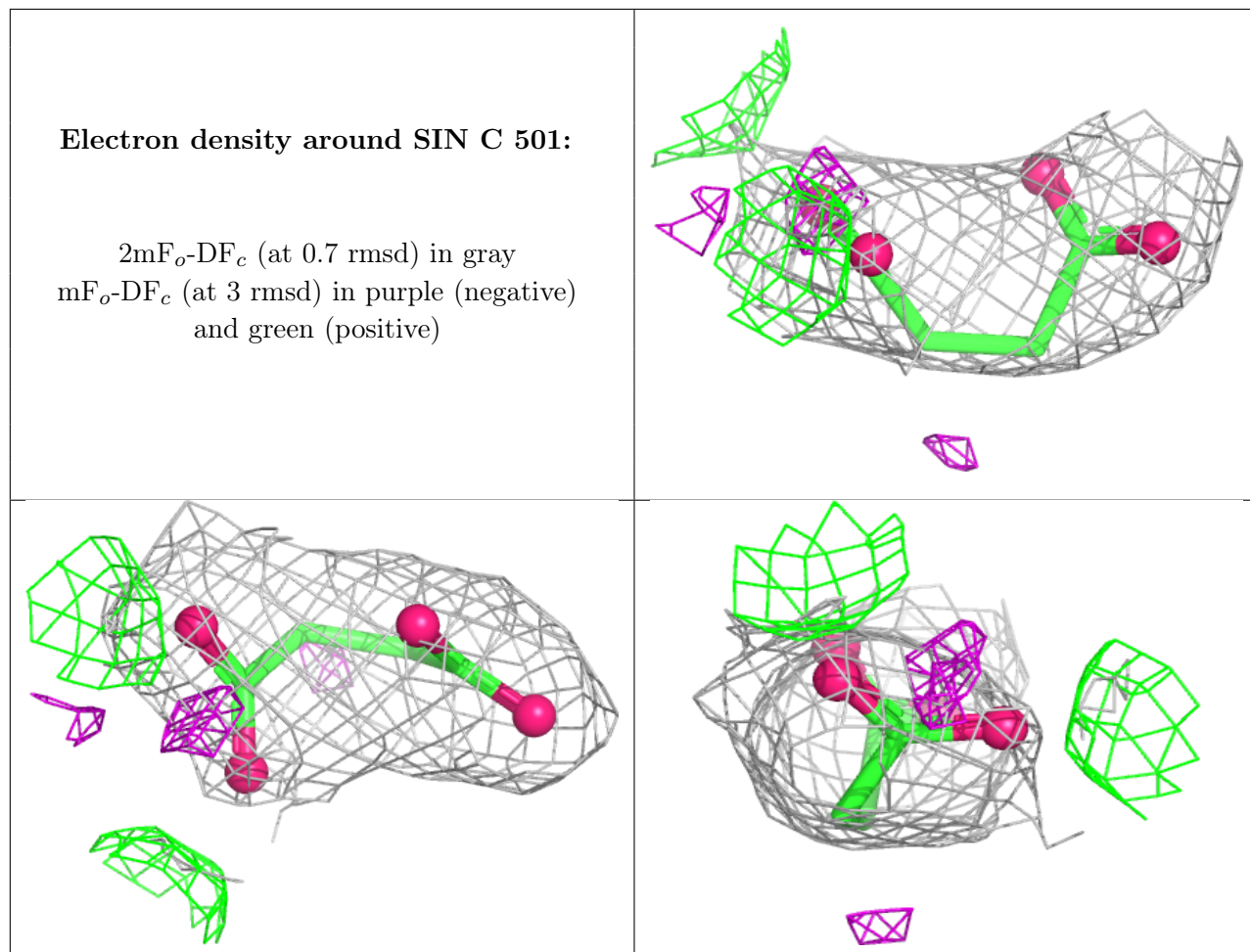
Electron density around SIN B 501:

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



Electron density around SIN C 501:

$2mF_o-DF_c$ (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.