



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 17, 2022 – 09:47 PM EDT

PDB ID : 3WZP
Title : Crystal structure of the core streptavidin mutant V21 (Y22S/N23D/S27D/Y83S/R84K/E101D/R103K/E116N) complexed with iminobiotin long tail (IMNtail) at 1.2 Å resolution
Authors : Kawato, T.; Mizohata, E.; Shimizu, Y.; Meshizuka, T.; Yamamoto, T.; Takasu, N.; Matsuoka, M.; Matsumura, H.; Tsumoto, K.; Kodama, T.; Kanai, M.; Doi, H.; Inoue, T.; Sugiyama, A.
Deposited on : 2014-10-01
Resolution : 1.20 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](#) ①) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.29
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) : 2.29

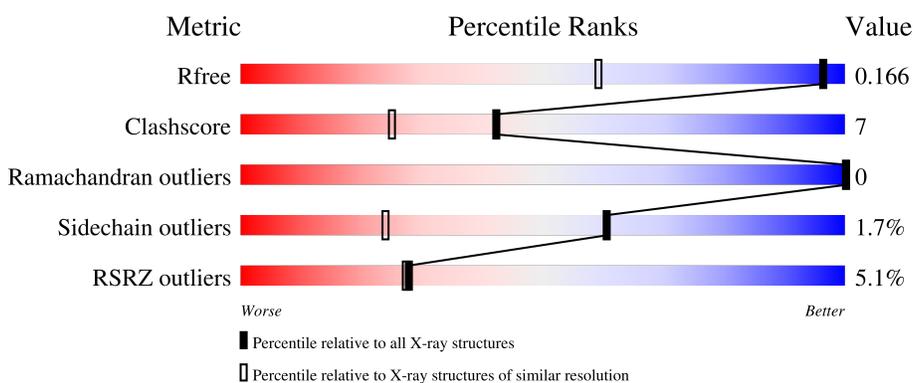
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.20 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	1223 (1.22-1.18)
Clashscore	141614	1286 (1.22-1.18)
Ramachandran outliers	138981	1240 (1.22-1.18)
Sidechain outliers	138945	1239 (1.22-1.18)
RSRZ outliers	127900	1200 (1.22-1.18)

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	129	<div style="display: flex; align-items: center;"> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 88%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">5% 88% 5% •• 5%</p>
1	B	129	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 84%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">3% 84% 10% • 5%</p>
1	C	129	<div style="display: flex; align-items: center;"> <div style="width: 9%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 83%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 9%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">9% 83% 9% •• 5%</p>
1	D	129	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 86%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 9%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">3% 86% 9% • 5%</p>

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
3	GOL	A	203	-	X	-	-
3	GOL	C	202	-	-	X	-
3	GOL	D	202	-	-	X	-
3	GOL	D	203	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4223 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 Streptavidin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
1	A	123	922	575	156	191	0	5	0
1	B	123	911	566	155	190	0	4	0
1	C	123	917	572	156	189	0	4	0
1	D	123	906	563	154	189	0	3	0

There are 40 discrepancies between the modelled and reference sequences:

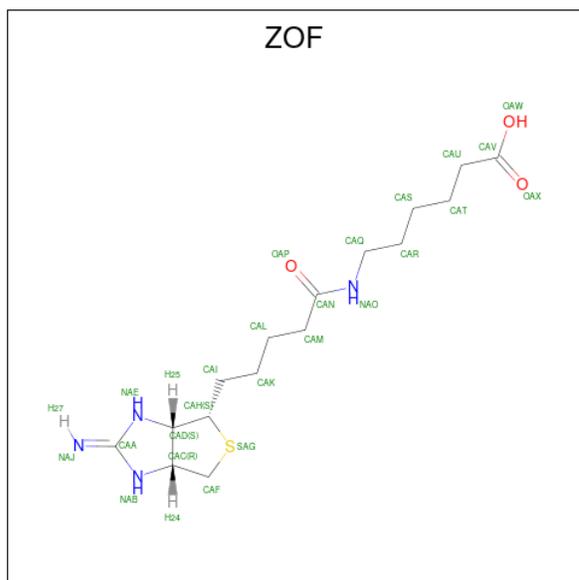
Chain	Residue	Modelled	Actual	Comment	Reference
A	11	GLY	-	expression tag	UNP P22629
A	12	PRO	-	expression tag	UNP P22629
A	22	SER	TYR	engineered mutation	UNP P22629
A	23	ASP	ASN	engineered mutation	UNP P22629
A	27	ASP	SER	engineered mutation	UNP P22629
A	83	SER	TYR	engineered mutation	UNP P22629
A	84	LYS	ARG	engineered mutation	UNP P22629
A	101	ASP	GLU	engineered mutation	UNP P22629
A	103	LYS	ARG	engineered mutation	UNP P22629
A	116	ASN	GLU	engineered mutation	UNP P22629
B	11	GLY	-	expression tag	UNP P22629
B	12	PRO	-	expression tag	UNP P22629
B	22	SER	TYR	engineered mutation	UNP P22629
B	23	ASP	ASN	engineered mutation	UNP P22629
B	27	ASP	SER	engineered mutation	UNP P22629
B	83	SER	TYR	engineered mutation	UNP P22629
B	84	LYS	ARG	engineered mutation	UNP P22629
B	101	ASP	GLU	engineered mutation	UNP P22629
B	103	LYS	ARG	engineered mutation	UNP P22629
B	116	ASN	GLU	engineered mutation	UNP P22629
C	11	GLY	-	expression tag	UNP P22629

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Chain	Residue	Modelled	Actual	Comment	Reference
C	12	PRO	-	expression tag	UNP P22629
C	22	SER	TYR	engineered mutation	UNP P22629
C	23	ASP	ASN	engineered mutation	UNP P22629
C	27	ASP	SER	engineered mutation	UNP P22629
C	83	SER	TYR	engineered mutation	UNP P22629
C	84	LYS	ARG	engineered mutation	UNP P22629
C	101	ASP	GLU	engineered mutation	UNP P22629
C	103	LYS	ARG	engineered mutation	UNP P22629
C	116	ASN	GLU	engineered mutation	UNP P22629
D	11	GLY	-	expression tag	UNP P22629
D	12	PRO	-	expression tag	UNP P22629
D	22	SER	TYR	engineered mutation	UNP P22629
D	23	ASP	ASN	engineered mutation	UNP P22629
D	27	ASP	SER	engineered mutation	UNP P22629
D	83	SER	TYR	engineered mutation	UNP P22629
D	84	LYS	ARG	engineered mutation	UNP P22629
D	101	ASP	GLU	engineered mutation	UNP P22629
D	103	LYS	ARG	engineered mutation	UNP P22629
D	116	ASN	GLU	engineered mutation	UNP P22629

- Molecule 2 is 6-({5-[(2E,3aS,4S,6aR)-2-iminohexahydro-1H-thieno[3,4-d]imidazol-4-yl]penta noyl}amino)hexanoic acid (three-letter code: ZOF) (formula: C₁₆H₂₈N₄O₃S).



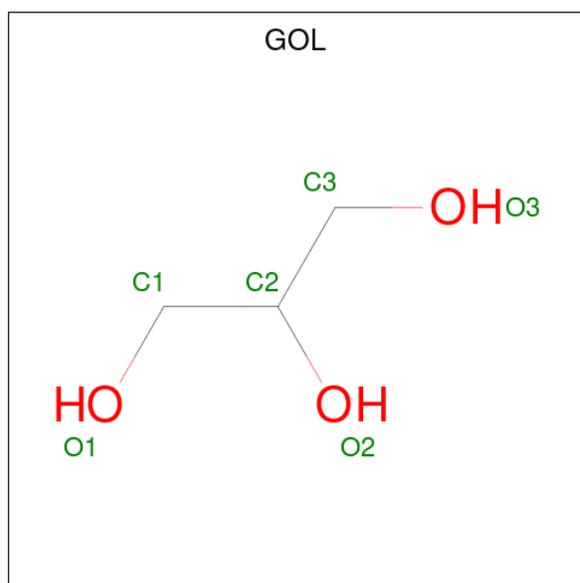
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
2	A	1	24	16	4	3	1	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	S	0	0
			24	16	4	3	1		
2	C	1	Total	C	N	O	S	0	0
			24	16	4	3	1		
2	D	1	Total	C	N	O	S	0	0
			24	16	4	3	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	125	Total	O	0	0
			125	125		

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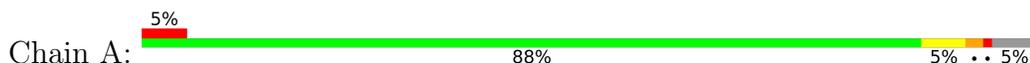
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	101	Total 101	O 101	0	0
4	C	110	Total 110	O 110	0	0
4	D	105	Total 105	O 105	0	0

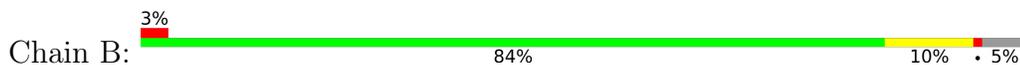
3 Residue-property plots [i](#)

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.

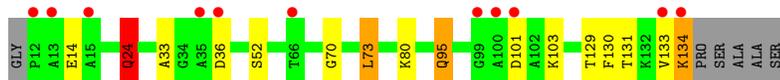
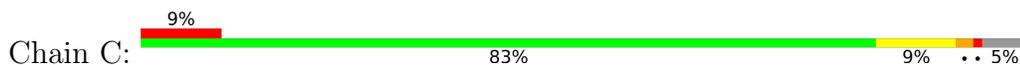
- Molecule 1: Streptavidin



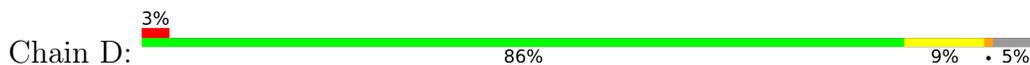
- Molecule 1: Streptavidin



- Molecule 1: Streptavidin



- Molecule 1: Streptavidin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.97Å 57.50Å 71.81Å 90.00° 103.34° 90.00°	Depositor
Resolution (Å)	15.00 – 1.20 15.00 – 1.20	Depositor EDS
% Data completeness (in resolution range)	96.5 (15.00-1.20) 96.5 (15.00-1.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.78 (at 1.20Å)	Xtrriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.134 , 0.161 0.142 , 0.166	Depositor DCC
R_{free} test set	7184 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	9.2	Xtrriage
Anisotropy	0.081	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.45 , 57.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4223	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.74% 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: ZOF, GOL

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	1.25	3/958 (0.3%)	1.15	2/1309 (0.2%)
1	B	1.21	7/944 (0.7%)	1.11	3/1293 (0.2%)
1	C	1.21	1/950 (0.1%)	1.14	2/1298 (0.2%)
1	D	1.27	4/936 (0.4%)	1.09	3/1282 (0.2%)
All	All	1.23	15/3788 (0.4%)	1.12	10/5182 (0.2%)

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	C	0	1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	51	GLU	CB-CG	-6.87	1.39	1.52
1	D	121	LYS	CB-CG	-6.71	1.34	1.52
1	C	95	GLN	CG-CD	6.40	1.65	1.51
1	A	95	GLN	CG-CD	5.86	1.64	1.51
1	B	116	ASN	CB-CG	5.85	1.64	1.51
1	B	24	GLN	CD-OE1	5.63	1.36	1.24
1	D	14	GLU	CD-OE2	5.50	1.31	1.25
1	B	51	GLU	CD-OE2	5.32	1.31	1.25
1	A	24	GLN	CB-CG	5.31	1.66	1.52
1	D	73	LEU	CG-CD2	-5.15	1.32	1.51
1	A	132	LYS	CB-CG	-5.15	1.38	1.52
1	B	24	GLN	CG-CD	5.13	1.62	1.51
1	B	73	LEU	CG-CD2	-5.09	1.33	1.51
1	B	121	LYS	CE-NZ	5.08	1.61	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	44	GLU	CD-OE1	5.03	1.31	1.25

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	73	LEU	CB-CG-CD2	7.95	124.51	111.00
1	C	73	LEU	CB-CG-CD1	7.93	124.47	111.00
1	B	36	ASP	CB-CG-OD1	7.41	124.97	118.30
1	D	27	ASP	CB-CG-OD1	6.51	124.16	118.30
1	B	27	ASP	CB-CG-OD1	6.17	123.85	118.30
1	D	36	ASP	CB-CG-OD2	6.08	123.77	118.30
1	D	27	ASP	CB-CG-OD2	-5.40	113.44	118.30
1	A	134	LYS	CD-CE-NZ	-5.29	99.53	111.70
1	C	133	VAL	CB-CA-C	-5.10	101.72	111.40
1	B	23	ASP	CB-CG-OD1	5.03	122.83	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	24	GLN	Sidechain

5.2 Too-close contacts

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	922	0	881	8	0
1	B	911	0	859	13	0
1	C	917	0	877	25	0
1	D	906	0	853	8	0
2	A	24	0	26	0	0
2	B	24	0	26	0	0
2	C	24	0	27	0	0
2	D	24	0	26	0	0
3	A	12	0	16	0	0
3	C	6	0	7	4	0
3	D	12	0	15	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	125	0	0	2	0
4	B	101	0	0	2	0
4	C	110	0	0	1	0
4	D	105	0	0	1	0
All	All	4223	0	3613	51	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 7.

All (51) 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:C:134:LYS:C	1:C:134:LYS:NZ	1.88	1.26
1:C:134:LYS:C	1:C:134:LYS:HZ2	1.45	1.15
1:C:36:ASP:O	3:C:202:GOL:O3	1.79	0.98
1:A:134:LYS:C	4:A:412:HOH:O	2.00	0.97
1:C:134:LYS:C	1:C:134:LYS:HZ3	1.65	0.93
1:A:95:GLN:HE22	1:B:114:THR:H	1.16	0.89
1:C:95:GLN:HE22	1:D:114:THR:H	1.18	0.89
1:B:49[A]:ASN:OD1	1:B:84:LYS:HE3	1.72	0.88
1:C:80:LYS:NZ	3:D:202:GOL:H12	1.90	0.87
1:B:49[A]:ASN:OD1	1:B:84:LYS:CE	2.27	0.83
1:C:80:LYS:HZ3	3:D:202:GOL:H12	1.50	0.76
1:D:100:ALA:O	1:D:101:ASP:HB3	1.94	0.67
1:C:103:LYS:NZ	1:C:131:THR:CG2	2.59	0.66
1:C:103:LYS:HD3	1:C:131:THR:HG22	1.77	0.65
1:C:52[A]:SER:OG	4:C:369:HOH:O	2.14	0.64
1:D:101:ASP:OD1	1:D:103:LYS:HE3	1.98	0.64
1:C:103:LYS:NZ	1:C:131:THR:HG21	2.14	0.63
1:C:103:LYS:CD	1:C:131:THR:HG22	2.28	0.62
1:A:95:GLN:NE2	1:B:114:THR:H	1.95	0.60
1:C:80:LYS:HZ2	3:D:202:GOL:H12	1.63	0.60
1:B:131:THR:HG23	4:B:346:HOH:O	2.07	0.55
1:C:103:LYS:HZ2	1:C:131:THR:CG2	2.19	0.54
1:B:103:LYS:NZ	1:B:131:THR:CG2	2.71	0.54
1:B:47:VAL:HG22	4:B:301:HOH:O	2.09	0.53
1:C:103:LYS:HZ3	1:C:131:THR:HG21	1.72	0.53
1:A:14:GLU:HG3	1:A:33:ALA:O	2.10	0.52
1:C:103:LYS:NZ	1:C:131:THR:HG22	2.24	0.52
1:B:103:LYS:NZ	1:B:131:THR:HG22	2.25	0.51
1:C:24:GLN:NE2	1:C:129:THR:H	2.08	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:36:ASP:O	3:D:202:GOL:O1	2.30	0.50
3:C:202:GOL:H31	1:D:80:LYS:HZ3	1.77	0.49
1:C:80:LYS:HZ2	3:D:202:GOL:C1	2.25	0.49
3:C:202:GOL:H31	1:D:80:LYS:NZ	2.27	0.49
1:A:70:GLY:HA3	1:A:95:GLN:HE21	1.78	0.48
1:C:103:LYS:HZ3	1:C:131:THR:CG2	2.27	0.47
1:D:116:ASN:HB3	4:D:349:HOH:O	2.13	0.47
1:B:24:GLN:CD	1:B:24:GLN:H	2.18	0.46
1:C:24:GLN:NE2	1:C:24:GLN:H	2.13	0.46
1:A:52[B]:SER:OG	4:A:363:HOH:O	2.21	0.46
1:C:14:GLU:HG3	1:C:33:ALA:O	2.16	0.45
1:C:70:GLY:HA3	1:C:95:GLN:HE21	1.81	0.45
1:B:103:LYS:HD3	1:B:131:THR:HG22	1.99	0.45
1:B:103:LYS:HZ2	1:B:131:THR:CG2	2.31	0.43
1:A:24:GLN:HE21	1:A:25:LEU:HG	1.83	0.42
3:C:202:GOL:C3	1:D:80:LYS:NZ	2.83	0.42
1:C:24:GLN:H	1:C:24:GLN:CD	2.24	0.41
1:C:101:ASP:O	1:C:103:LYS:HG2	2.20	0.41
1:B:103:LYS:NZ	1:B:131:THR:HG21	2.34	0.41
1:B:103:LYS:HZ3	1:B:131:THR:CG2	2.33	0.41
1:A:24:GLN:CD	1:A:24:GLN:H	2.24	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	126/129 (98%)	121 (96%)	5 (4%)	0	100	100
1	B	125/129 (97%)	120 (96%)	5 (4%)	0	100	100
1	C	125/129 (97%)	121 (97%)	4 (3%)	0	100	100
1	D	124/129 (96%)	119 (96%)	5 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	500/516 (97%)	481 (96%)	19 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	95/93 (102%)	93 (98%)	2 (2%)	53	15
1	B	93/93 (100%)	92 (99%)	1 (1%)	73	41
1	C	94/93 (101%)	91 (97%)	3 (3%)	39	6
1	D	92/93 (99%)	92 (100%)	0	100	100
All	All	374/372 (100%)	368 (98%)	6 (2%)	60	27

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	73	LEU
1	B	24	GLN
1	C	24	GLN
1	C	73	LEU
1	C	134	LYS

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

Mol	Chain	Res	Type
1	A	95	GLN
1	A	107	GLN
1	A	116	ASN
1	C	24	GLN
1	C	95	GLN
1	C	107	GLN

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Mol	Chain	Res	Type
1	D	116	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

9 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	ZOF	C	201	-	25,25,25	1.18	5 (20%)	29,32,32	1.54	6 (20%)
3	GOL	C	202	-	5,5,5	0.84	0	5,5,5	1.82	1 (20%)
3	GOL	A	202	-	5,5,5	0.53	0	5,5,5	3.78	2 (40%)
2	ZOF	A	201	-	25,25,25	1.03	2 (8%)	29,32,32	1.80	8 (27%)
3	GOL	D	203	-	5,5,5	0.70	0	5,5,5	4.18	3 (60%)
2	ZOF	B	201	-	25,25,25	2.03	7 (28%)	29,32,32	2.35	9 (31%)
3	GOL	A	203	-	5,5,5	0.77	0	5,5,5	2.20	2 (40%)
2	ZOF	D	201	-	25,25,25	1.83	8 (32%)	29,32,32	2.00	8 (27%)
3	GOL	D	202	-	5,5,5	1.19	0	5,5,5	2.09	1 (20%)

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	ZOF	C	201	-	-	2/16/37/37	0/2/2/2
3	GOL	C	202	-	-	2/4/4/4	-
3	GOL	A	202	-	-	3/4/4/4	-
2	ZOF	A	201	-	-	2/16/37/37	0/2/2/2
3	GOL	D	203	-	-	3/4/4/4	-
2	ZOF	B	201	-	-	2/16/37/37	0/2/2/2
3	GOL	A	203	-	-	4/4/4/4	-
2	ZOF	D	201	-	-	2/16/37/37	0/2/2/2
3	GOL	D	202	-	-	0/4/4/4	-

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	201	ZOF	CAF-CAC	5.58	1.63	1.53
2	B	201	ZOF	CAH-CAD	4.33	1.63	1.53
2	D	201	ZOF	CAF-CAC	3.95	1.60	1.53
2	D	201	ZOF	CAA-NAB	-3.72	1.29	1.35
2	B	201	ZOF	CAD-NAE	3.53	1.52	1.45
2	B	201	ZOF	CAH-SAG	-3.29	1.77	1.82
2	D	201	ZOF	CAH-CAD	2.99	1.60	1.53
2	D	201	ZOF	CAU-CAV	2.98	1.57	1.50
2	D	201	ZOF	CAH-SAG	-2.87	1.78	1.82
2	D	201	ZOF	OAX-CAV	2.70	1.31	1.22
2	B	201	ZOF	CAU-CAV	2.55	1.56	1.50
2	A	201	ZOF	CAU-CAV	2.51	1.56	1.50
2	C	201	ZOF	CAA-NAE	-2.44	1.31	1.35
2	D	201	ZOF	CAF-SAG	-2.36	1.74	1.81
2	C	201	ZOF	CAA-NAB	-2.28	1.31	1.35
2	B	201	ZOF	CAA-NAB	-2.27	1.31	1.35
2	D	201	ZOF	CAM-CAN	2.13	1.55	1.51
2	C	201	ZOF	CAU-CAV	2.13	1.55	1.50
2	C	201	ZOF	CAA-NAJ	2.11	1.35	1.29
2	B	201	ZOF	OAX-CAV	2.07	1.29	1.22
2	A	201	ZOF	OAX-CAV	2.07	1.29	1.22
2	C	201	ZOF	OAX-CAV	2.04	1.28	1.22

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	203	GOL	C3-C2-C1	-8.75	77.70	111.70
3	A	202	GOL	C3-C2-C1	-7.43	82.81	111.70
2	D	201	ZOF	NAE-CAA-NAB	5.45	114.80	109.20
2	B	201	ZOF	CAD-CAH-SAG	-5.19	100.26	105.20
2	B	201	ZOF	CAF-CAC-CAD	-5.10	104.23	108.66
2	B	201	ZOF	NAE-CAA-NAB	4.78	114.10	109.20
2	A	201	ZOF	CAF-CAC-NAB	-4.55	107.25	113.03
2	A	201	ZOF	NAE-CAA-NAB	4.42	113.74	109.20
2	B	201	ZOF	CAF-CAC-NAB	-4.21	107.68	113.03
2	C	201	ZOF	NAE-CAA-NAB	3.87	113.17	109.20
3	A	203	GOL	O3-C3-C2	-3.83	91.84	110.20
2	D	201	ZOF	CAF-CAC-CAD	-3.78	105.38	108.66
2	C	201	ZOF	CAH-CAD-NAE	3.64	116.39	113.13
3	D	202	GOL	C3-C2-C1	-3.61	97.66	111.70
2	B	201	ZOF	CAK-CAI-CAH	-3.40	106.96	113.86
2	D	201	ZOF	CAD-CAH-SAG	-3.37	101.99	105.20
2	B	201	ZOF	CAH-CAD-NAE	3.37	116.14	113.13
2	D	201	ZOF	CAH-CAD-NAE	3.29	116.08	113.13
2	B	201	ZOF	CAD-NAE-CAA	-3.27	106.65	112.08
3	C	202	GOL	O2-C2-C3	3.03	122.45	109.12
3	A	202	GOL	O2-C2-C1	3.02	122.44	109.12
2	C	201	ZOF	CAF-CAC-NAB	-2.93	109.30	113.03
2	B	201	ZOF	CAC-CAD-NAE	2.88	105.77	102.67
2	D	201	ZOF	CAT-CAU-CAV	-2.70	107.67	114.47
2	A	201	ZOF	CAK-CAI-CAH	-2.60	108.58	113.86
2	D	201	ZOF	CAM-CAN-NAO	-2.57	112.10	116.42
2	C	201	ZOF	CAD-CAH-SAG	-2.56	102.76	105.20
2	D	201	ZOF	CAK-CAI-CAH	-2.54	108.71	113.86
2	A	201	ZOF	CAH-CAD-CAC	-2.45	106.10	108.94
2	A	201	ZOF	CAF-CAC-CAD	2.44	110.78	108.66
2	A	201	ZOF	CAH-CAD-NAE	2.41	115.28	113.13
2	C	201	ZOF	CAF-CAC-CAD	-2.34	106.63	108.66
2	A	201	ZOF	CAD-CAC-NAB	2.33	104.91	102.43
2	D	201	ZOF	CAS-CAT-CAU	-2.32	104.86	113.19
2	A	201	ZOF	CAD-CAH-SAG	-2.31	103.00	105.20
2	C	201	ZOF	CAK-CAI-CAH	-2.25	109.29	113.86
3	D	203	GOL	O3-C3-C2	-2.17	99.81	110.20
3	D	203	GOL	O2-C2-C1	2.15	118.59	109.12
2	B	201	ZOF	CAT-CAU-CAV	-2.06	109.29	114.47
3	A	203	GOL	C3-C2-C1	-2.03	103.83	111.70

There are no chirality outliers.

All (20) torsion outliers are listed below:

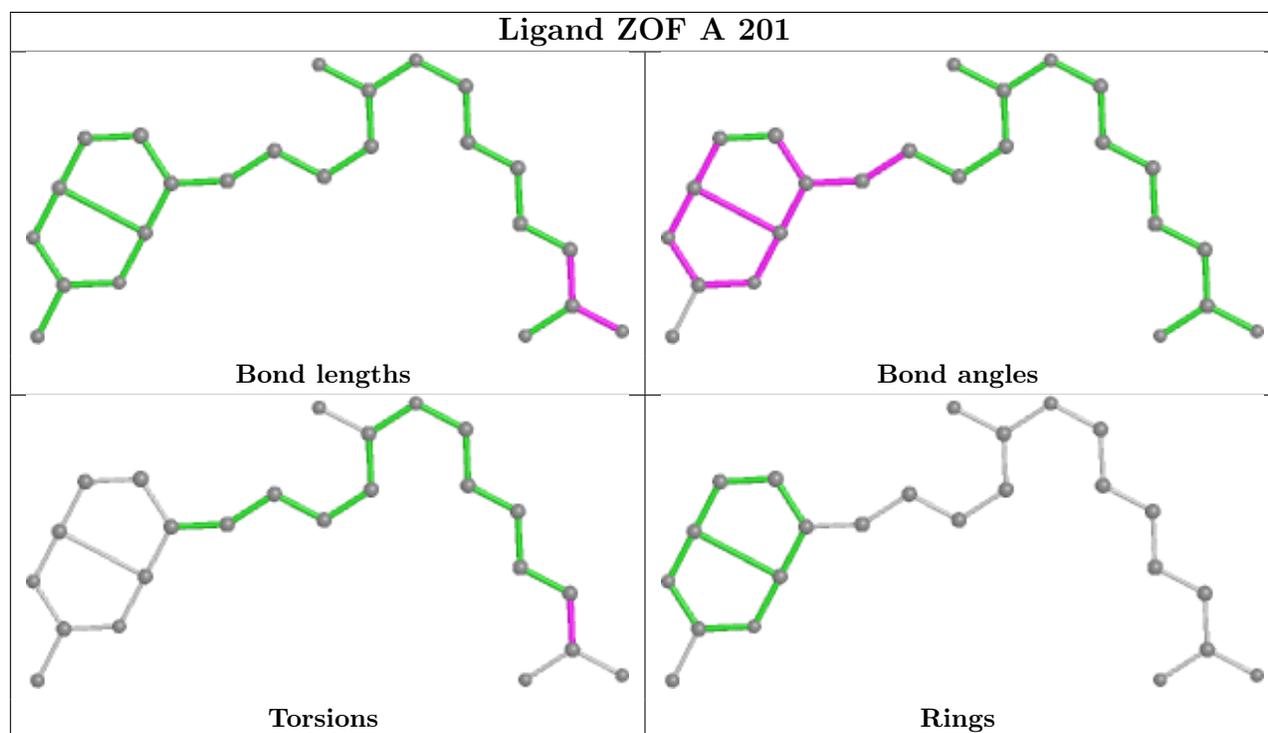
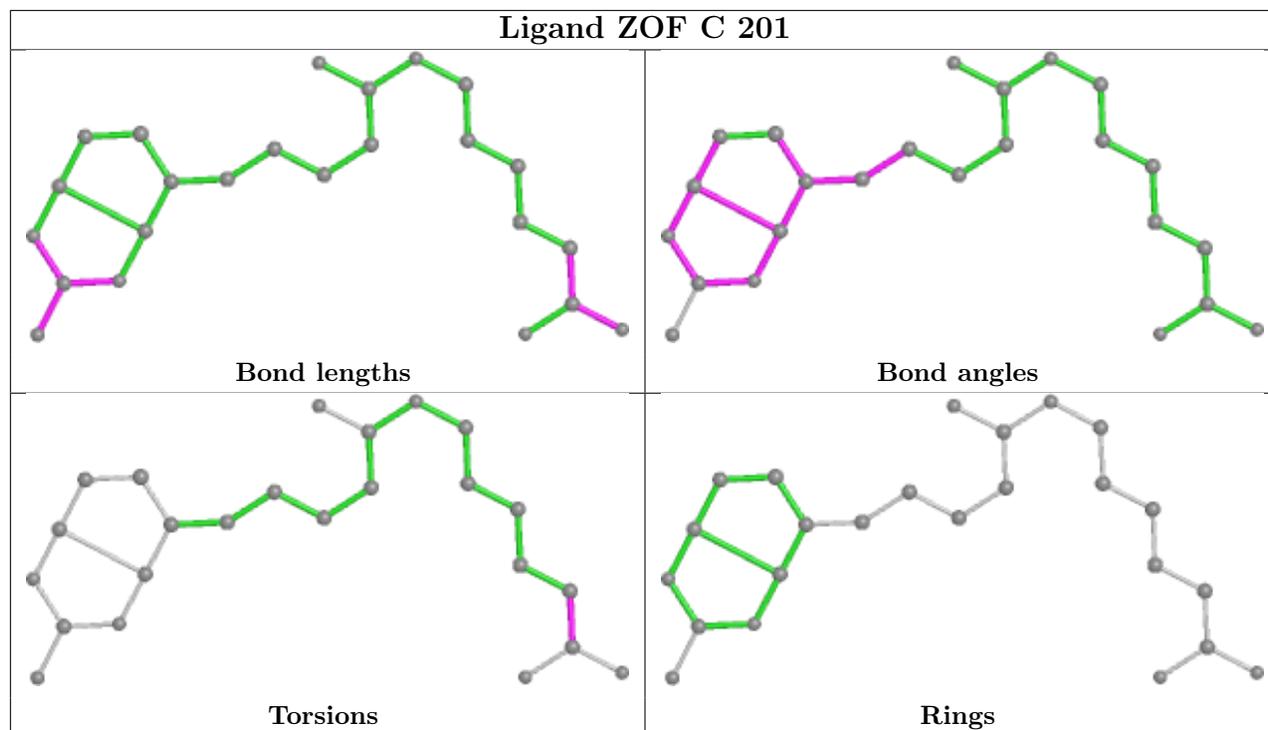
Mol	Chain	Res	Type	Atoms
3	A	202	GOL	C1-C2-C3-O3
3	A	203	GOL	O1-C1-C2-C3
3	C	202	GOL	C1-C2-C3-O3
3	D	203	GOL	C1-C2-C3-O3
3	D	203	GOL	O2-C2-C3-O3
3	A	203	GOL	O1-C1-C2-O2
3	A	203	GOL	O2-C2-C3-O3
3	A	202	GOL	O1-C1-C2-C3
3	A	203	GOL	C1-C2-C3-O3
3	D	203	GOL	O1-C1-C2-C3
3	A	202	GOL	O2-C2-C3-O3
3	C	202	GOL	O2-C2-C3-O3
2	D	201	ZOF	CAT-CAU-CAV-OAW
2	D	201	ZOF	CAT-CAU-CAV-OAX
2	B	201	ZOF	CAT-CAU-CAV-OAX
2	A	201	ZOF	CAT-CAU-CAV-OAX
2	B	201	ZOF	CAT-CAU-CAV-OAW
2	C	201	ZOF	CAT-CAU-CAV-OAW
2	C	201	ZOF	CAT-CAU-CAV-OAX
2	A	201	ZOF	CAT-CAU-CAV-OAW

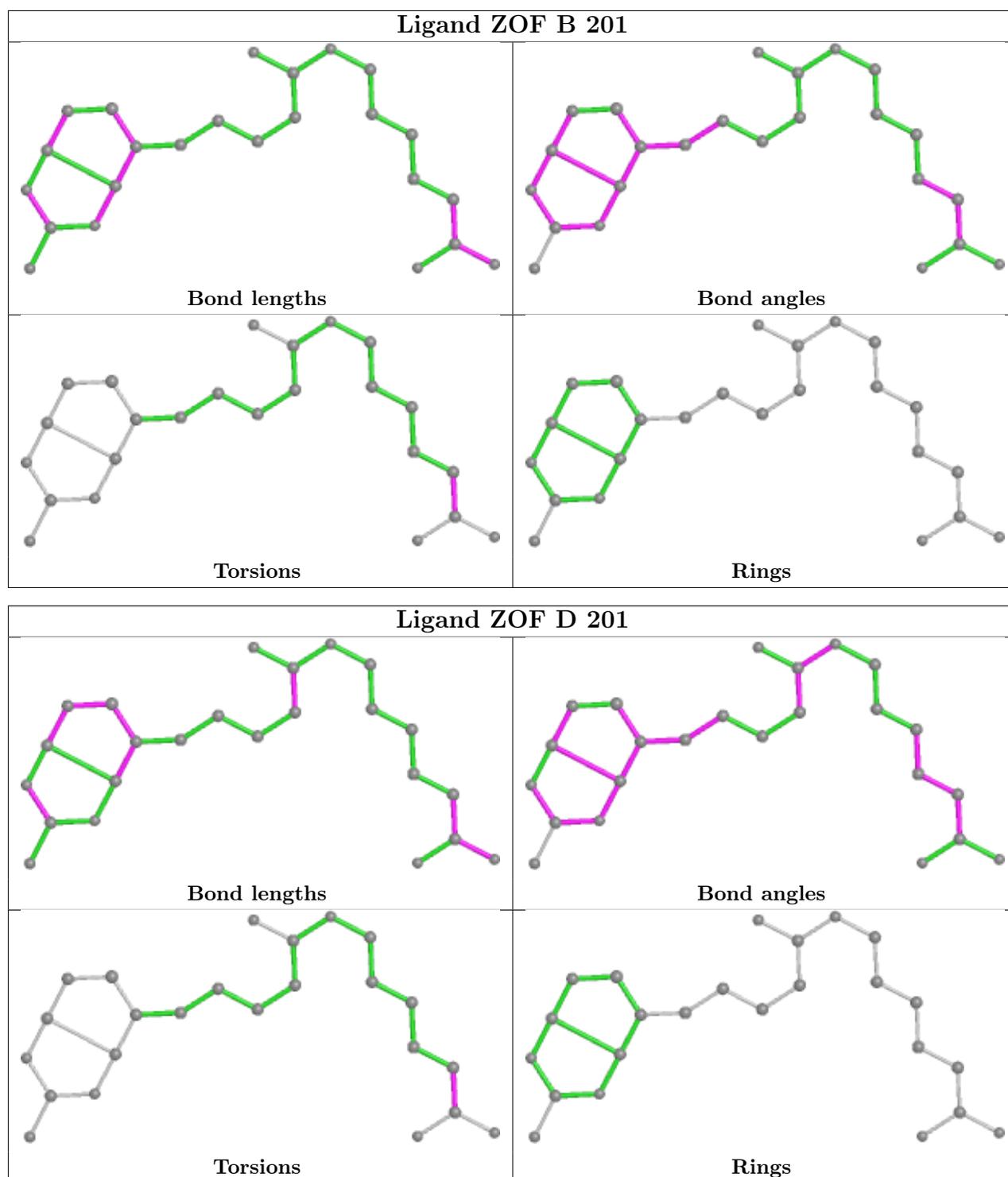
There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	202	GOL	4	0
3	D	202	GOL	5	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

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	123/129 (95%)	-0.10	6 (4%) 29 29	5, 7, 19, 37	0
1	B	123/129 (95%)	-0.16	4 (3%) 46 45	5, 8, 19, 38	0
1	C	123/129 (95%)	0.02	11 (8%) 9 10	5, 8, 24, 49	0
1	D	123/129 (95%)	-0.14	4 (3%) 46 45	5, 9, 20, 37	0
All	All	492/516 (95%)	-0.10	25 (5%) 28 27	5, 8, 22, 49	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	100	ALA	8.0
1	A	100	ALA	7.2
1	C	134	LYS	5.1
1	C	12	PRO	5.0
1	B	101	ASP	5.0
1	D	101	ASP	4.5
1	A	101	ASP	4.4
1	B	100	ALA	3.9
1	A	12	PRO	3.9
1	D	100	ALA	3.8
1	A	15	ALA	3.7
1	C	66	THR	3.6
1	A	99	GLY	3.5
1	C	15	ALA	3.3
1	C	101	ASP	3.2
1	A	134	LYS	3.1
1	C	35	ALA	2.9
1	D	133	VAL	2.8
1	C	99	GLY	2.6
1	C	133	VAL	2.4
1	C	36	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	13	ALA	2.3
1	B	133	VAL	2.2
1	D	24	GLN	2.1
1	B	24	GLN	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](#)

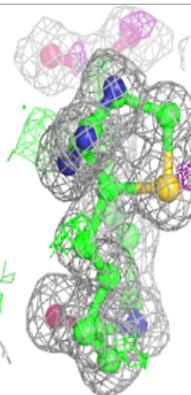
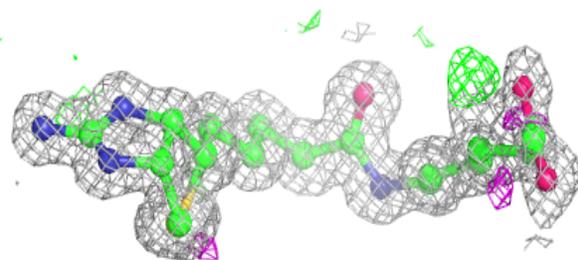
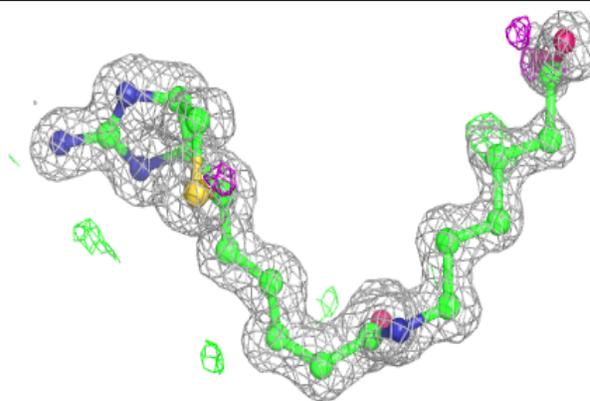
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GOL	C	202	6/6	0.70	0.29	18,23,28,29	6
3	GOL	A	203	6/6	0.83	0.25	17,18,23,27	6
3	GOL	D	202	6/6	0.86	0.25	18,22,24,25	6
3	GOL	D	203	6/6	0.88	0.35	10,15,19,21	6
3	GOL	A	202	6/6	0.93	0.33	6,14,18,19	6
2	ZOF	D	201	24/24	0.97	0.09	8,12,30,37	0
2	ZOF	B	201	24/24	0.97	0.09	7,11,27,31	0
2	ZOF	A	201	24/24	0.98	0.07	5,7,20,25	0
2	ZOF	C	201	24/24	0.98	0.08	6,7,19,29	0

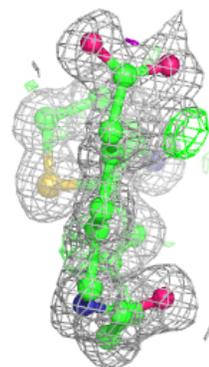
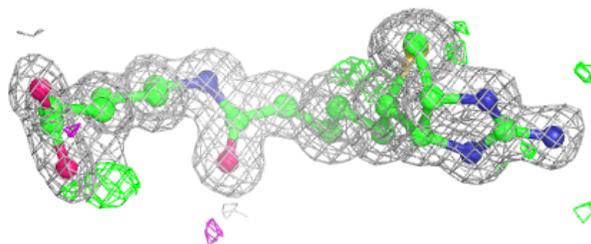
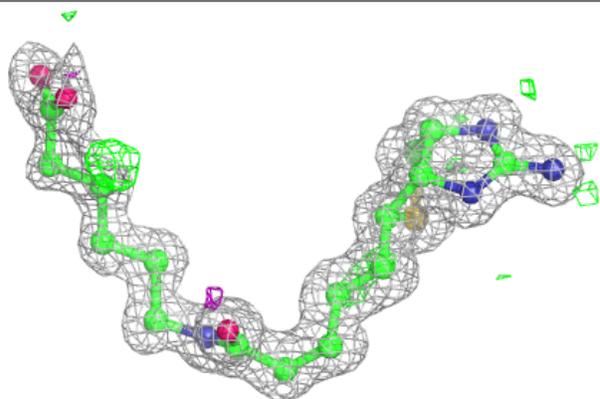
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ZOF D 201:

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

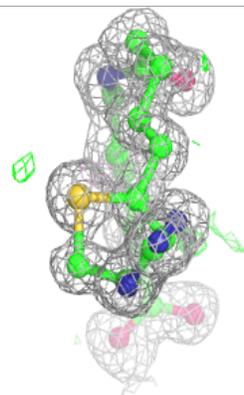
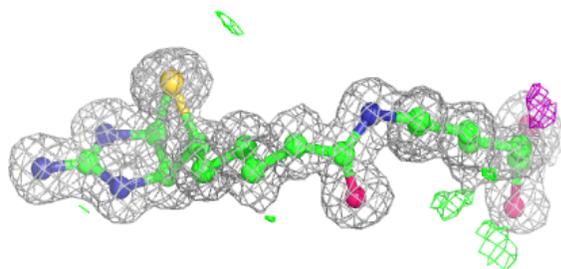
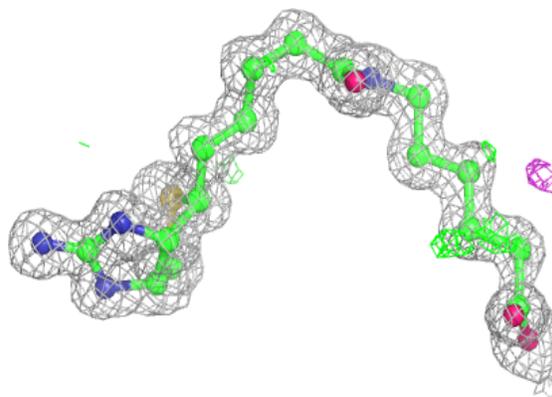
**Electron density around ZOF B 201:**

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

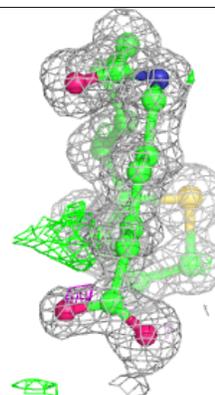
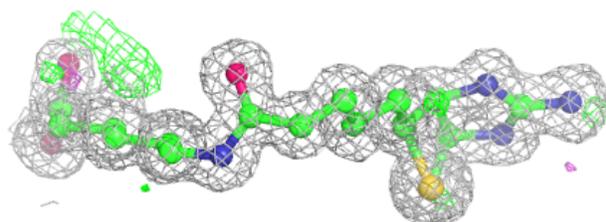
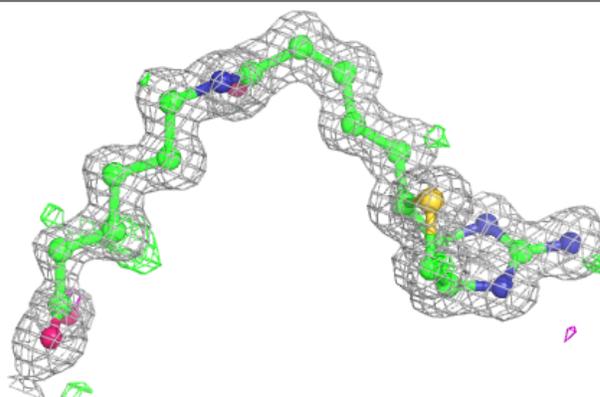


Electron density around ZOF A 201:

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

**Electron density around ZOF C 201:**

$2mF_o-DF_c$ (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.