



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 06:00 PM EDT

PDB ID : 3WZQ / pdb_00003wzq
Title : Crystal structure of the core streptavidin mutant V212 (Y22S/N23D/S27D/S45N/Y83S/R84K/E101D/R103K/E116N) complexed with iminobiotin long tail (IMNtail) at 1.7 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 : 1.70 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)

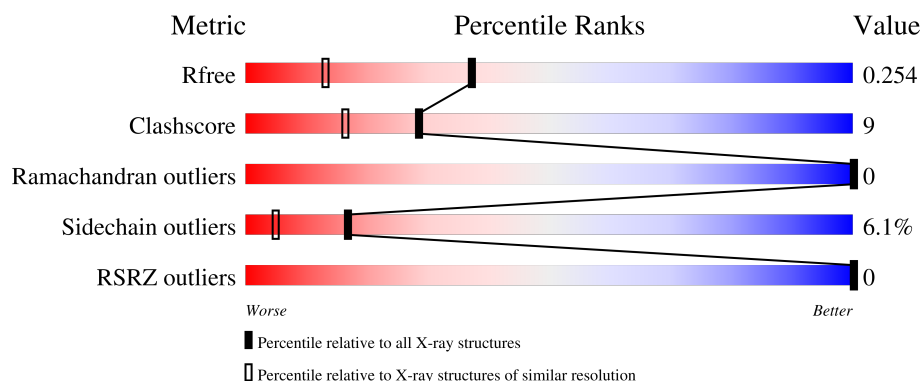
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	147	
1	B	147	
1	C	147	

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Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

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Mol	Chain	Length	Quality of chain
1	D	147	 69% 13% •• 15%

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	P6G	D	1002	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4128 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
1	A	125	Total	C	N	O	0	6	0
			945	593	158	194			
1	B	125	Total	C	N	O	0	5	0
			941	590	158	193			
1	C	124	Total	C	N	O	0	1	0
			910	564	157	189			
1	D	125	Total	C	N	O	0	1	0
			917	569	158	190			

There are 116 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	expression tag	UNP P22629
A	0	ALA	-	expression tag	UNP P22629
A	1	SER	-	expression tag	UNP P22629
A	2	MET	-	expression tag	UNP P22629
A	3	THR	-	expression tag	UNP P22629
A	4	GLY	-	expression tag	UNP P22629
A	5	GLY	-	expression tag	UNP P22629
A	6	GLN	-	expression tag	UNP P22629
A	7	GLN	-	expression tag	UNP P22629
A	8	MET	-	expression tag	UNP P22629
A	9	GLY	-	expression tag	UNP P22629
A	10	ARG	-	expression tag	UNP P22629
A	11	GLY	-	expression tag	UNP P22629
A	12	SER	-	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	45	ASN	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

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Chain	Residue	Modelled	Actual	Comment	Reference
A	103	LYS	ARG	engineered mutation	UNP P22629
A	116	ASN	GLU	engineered mutation	UNP P22629
A	140	HIS	-	expression tag	UNP P22629
A	141	HIS	-	expression tag	UNP P22629
A	142	HIS	-	expression tag	UNP P22629
A	143	HIS	-	expression tag	UNP P22629
A	144	HIS	-	expression tag	UNP P22629
A	145	HIS	-	expression tag	UNP P22629
B	-1	MET	-	expression tag	UNP P22629
B	0	ALA	-	expression tag	UNP P22629
B	1	SER	-	expression tag	UNP P22629
B	2	MET	-	expression tag	UNP P22629
B	3	THR	-	expression tag	UNP P22629
B	4	GLY	-	expression tag	UNP P22629
B	5	GLY	-	expression tag	UNP P22629
B	6	GLN	-	expression tag	UNP P22629
B	7	GLN	-	expression tag	UNP P22629
B	8	MET	-	expression tag	UNP P22629
B	9	GLY	-	expression tag	UNP P22629
B	10	ARG	-	expression tag	UNP P22629
B	11	GLY	-	expression tag	UNP P22629
B	12	SER	-	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	45	ASN	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
B	140	HIS	-	expression tag	UNP P22629
B	141	HIS	-	expression tag	UNP P22629
B	142	HIS	-	expression tag	UNP P22629
B	143	HIS	-	expression tag	UNP P22629
B	144	HIS	-	expression tag	UNP P22629
B	145	HIS	-	expression tag	UNP P22629
C	-1	MET	-	expression tag	UNP P22629
C	0	ALA	-	expression tag	UNP P22629
C	1	SER	-	expression tag	UNP P22629
C	2	MET	-	expression tag	UNP P22629
C	3	THR	-	expression tag	UNP P22629

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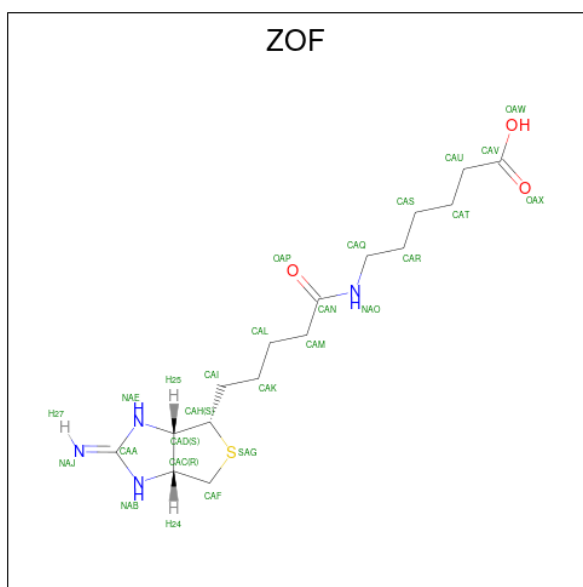
Chain	Residue	Modelled	Actual	Comment	Reference
C	4	GLY	-	expression tag	UNP P22629
C	5	GLY	-	expression tag	UNP P22629
C	6	GLN	-	expression tag	UNP P22629
C	7	GLN	-	expression tag	UNP P22629
C	8	MET	-	expression tag	UNP P22629
C	9	GLY	-	expression tag	UNP P22629
C	10	ARG	-	expression tag	UNP P22629
C	11	GLY	-	expression tag	UNP P22629
C	12	SER	-	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	45	ASN	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
C	140	HIS	-	expression tag	UNP P22629
C	141	HIS	-	expression tag	UNP P22629
C	142	HIS	-	expression tag	UNP P22629
C	143	HIS	-	expression tag	UNP P22629
C	144	HIS	-	expression tag	UNP P22629
C	145	HIS	-	expression tag	UNP P22629
D	-1	MET	-	expression tag	UNP P22629
D	0	ALA	-	expression tag	UNP P22629
D	1	SER	-	expression tag	UNP P22629
D	2	MET	-	expression tag	UNP P22629
D	3	THR	-	expression tag	UNP P22629
D	4	GLY	-	expression tag	UNP P22629
D	5	GLY	-	expression tag	UNP P22629
D	6	GLN	-	expression tag	UNP P22629
D	7	GLN	-	expression tag	UNP P22629
D	8	MET	-	expression tag	UNP P22629
D	9	GLY	-	expression tag	UNP P22629
D	10	ARG	-	expression tag	UNP P22629
D	11	GLY	-	expression tag	UNP P22629
D	12	SER	-	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	45	ASN	SER	engineered mutation	UNP P22629

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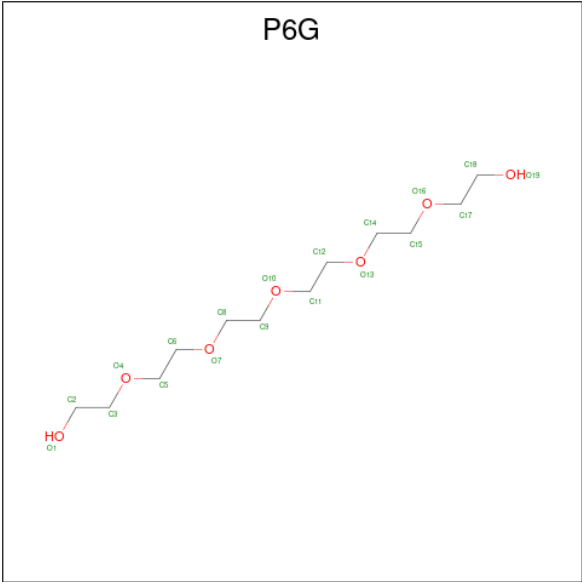
Chain	Residue	Modelled	Actual	Comment	Reference
D	83	SER	TYR	engineered mutation	UNP P22629
D	84	LYS	ARG	engineered mutation	UNP P22629
D	101	ASP	GLU	engineered mutation	UNP P22629
D	103	LYS	ARG	engineered mutation	UNP P22629
D	116	ASN	GLU	engineered mutation	UNP P22629
D	140	HIS	-	expression tag	UNP P22629
D	141	HIS	-	expression tag	UNP P22629
D	142	HIS	-	expression tag	UNP P22629
D	143	HIS	-	expression tag	UNP P22629
D	144	HIS	-	expression tag	UNP P22629
D	145	HIS	-	expression tag	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 (CCD ID: ZOF) (formula: C₁₆H₂₈N₄O₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			24	16	4	3	1		
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 HEXAETHYLENE GLYCOL (CCD ID: P6G) (formula: C₁₂H₂₆O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	C	O	0	0
			16	10	6		
3	D	1	Total	C	O	0	0
			16	10	6		

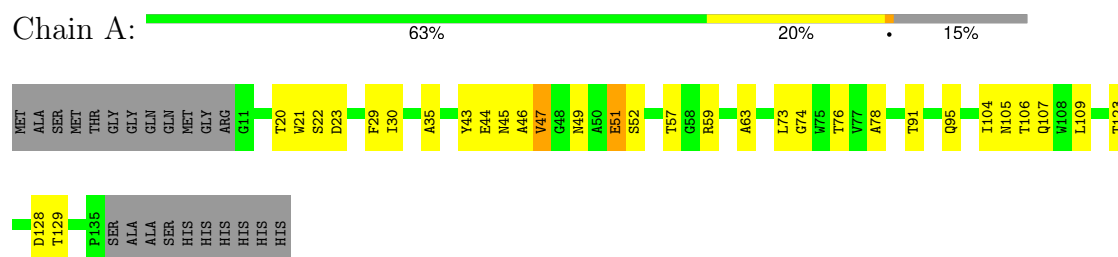
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	81	Total	O	0	0
			81	81		
4	B	83	Total	O	0	0
			83	83		
4	C	62	Total	O	0	0
			62	62		
4	D	61	Total	O	0	0
			61	61		

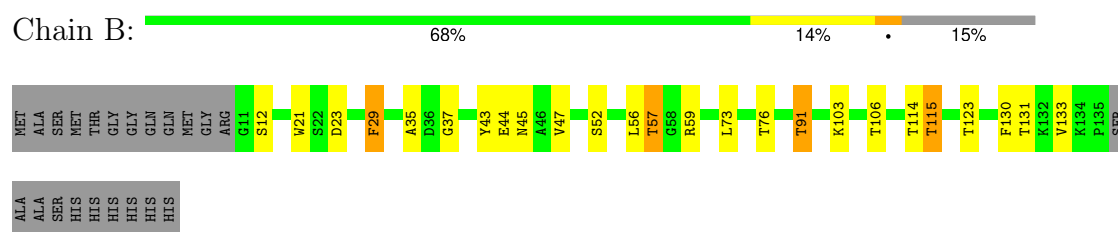
3 Residue-property plots [i](#)

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.

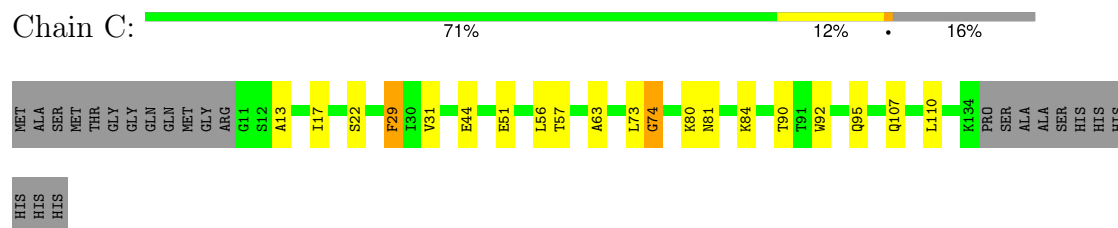
• Molecule 1: Streptavidin



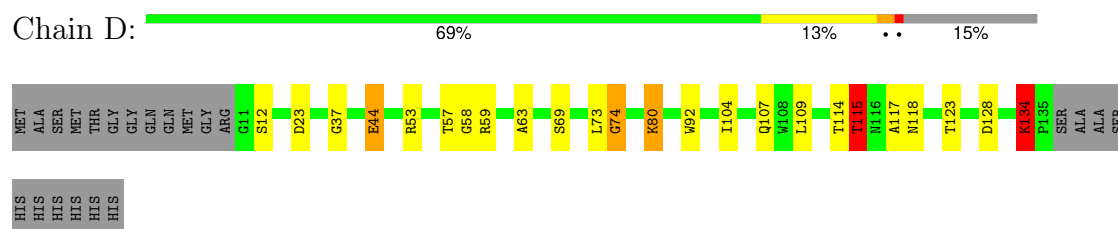
• Molecule 1: Streptavidin



• Molecule 1: Streptavidin



• Molecule 1: Streptavidin



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	77.42Å 77.41Å 172.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 1.70 15.00 – 1.70	Depositor EDS
% Data completeness (in resolution range)	96.8 (15.00-1.70) 96.8 (15.00-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.24 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.197 , 0.244 0.207 , 0.254	Depositor DCC
R_{free} test set	2812 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	17.3	Xtriage
Anisotropy	0.244	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 35.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.477 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4128	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.47% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZOF, P6G

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	1.50	9/986 (0.9%)	1.20	4/1350 (0.3%)
1	B	1.50	5/979 (0.5%)	1.23	3/1340 (0.2%)
1	C	1.42	1/933 (0.1%)	1.23	3/1276 (0.2%)
1	D	1.49	6/941 (0.6%)	1.26	7/1288 (0.5%)
All	All	1.48	21/3839 (0.5%)	1.23	17/5254 (0.3%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	74	GLY	N-CA	7.40	1.52	1.45
1	D	37	GLY	N-CA	6.88	1.54	1.45
1	A	35	ALA	CA-CB	6.67	1.65	1.53
1	B	35	ALA	CA-CB	6.37	1.64	1.53
1	D	80	LYS	N-CA	-6.34	1.38	1.46
1	D	58	GLY	C-O	6.13	1.30	1.23
1	D	74	GLY	N-CA	6.08	1.51	1.45
1	A	106	THR	N-CA	5.80	1.53	1.45
1	B	37	GLY	N-CA	5.78	1.53	1.45
1	B	91	THR	N-CA	5.73	1.53	1.46
1	A	109	LEU	C-O	5.50	1.30	1.24
1	A	78	ALA	CA-CB	5.43	1.61	1.53
1	A	123	THR	N-CA	5.38	1.52	1.46
1	B	123	THR	N-CA	5.35	1.52	1.46
1	D	104	ILE	CA-CB	-5.34	1.47	1.54
1	A	104	ILE	CA-CB	-5.27	1.47	1.54
1	B	76	THR	C-O	5.23	1.30	1.23
1	A	76	THR	C-O	5.17	1.30	1.24
1	A	74	GLY	N-CA	5.14	1.50	1.45
1	A	107	GLN	C-O	5.09	1.30	1.23
1	D	134	LYS	CA-C	5.03	1.57	1.52

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	115	THR	N-CA-CB	-7.84	99.15	110.36
1	C	17	ILE	N-CA-C	7.61	117.73	110.42
1	B	106	THR	OG1-CB-CG2	6.93	123.15	109.30
1	D	69	SER	CA-C-N	-6.35	113.19	121.96
1	D	69	SER	C-N-CA	-6.35	113.19	121.96
1	D	123	THR	N-CA-C	6.05	118.26	108.34
1	A	63	ALA	CA-C-N	-5.95	114.61	120.98
1	A	63	ALA	C-N-CA	-5.95	114.61	120.98
1	B	12	SER	N-CA-C	5.90	117.52	111.14
1	C	107	GLN	CA-CB-CG	-5.77	102.56	114.10
1	A	107	GLN	CA-CB-CG	-5.70	102.70	114.10
1	D	107	GLN	CA-CB-CG	-5.69	102.72	114.10
1	C	13	ALA	N-CA-C	-5.66	105.20	111.71
1	A	46	ALA	N-CA-C	5.46	117.23	111.28
1	D	109	LEU	CA-C-N	-5.22	115.64	123.00
1	D	109	LEU	C-N-CA	-5.22	115.64	123.00
1	D	115	THR	N-CA-CB	-5.12	103.04	110.36

There are no chirality outliers.

There are no planarity outliers.

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	945	0	899	20	0
1	B	941	0	892	15	0
1	C	910	0	855	20	0
1	D	917	0	862	20	0
2	A	24	0	26	0	0
2	B	24	0	27	0	0
2	C	24	0	27	0	0
2	D	24	0	26	0	0
3	C	16	0	21	7	0
3	D	16	0	21	10	0
4	A	81	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	83	0	0	1	0
4	C	62	0	0	2	0
4	D	61	0	0	0	0
All	All	4128	0	3656	66	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 9.

All (66) 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:29:PHE:CE1	1:C:31:VAL:HB	1.94	1.03
1:C:29:PHE:HE1	1:C:31:VAL:HB	1.25	0.97
1:C:95:GLN:HE22	1:D:114:THR:H	1.19	0.89
1:D:134:LYS:HZ2	1:D:134:LYS:HA	1.42	0.83
1:A:20:THR:HG23	1:A:30:ILE:CD1	2.15	0.77
1:A:95:GLN:HE22	1:B:114:THR:H	1.32	0.77
1:C:29:PHE:HE1	1:C:31:VAL:CB	1.97	0.76
1:A:59:ARG:HD3	1:B:57[A]:THR:HG23	1.68	0.76
1:A:51:GLU:H	1:A:51:GLU:CD	1.90	0.75
1:A:45:ASN:HD21	1:A:47:VAL:HG23	1.56	0.71
1:D:80:LYS:HZ2	3:D:1002:P6G:H122	1.56	0.71
1:C:63:ALA:HB2	3:D:1002:P6G:H152	1.75	0.69
1:A:57[A]:THR:HG22	1:B:59:ARG:HG2	1.76	0.67
1:D:80:LYS:NZ	3:D:1002:P6G:H91	2.10	0.66
1:A:105:ASN:OD1	1:A:129[B]:THR:HG22	1.96	0.65
1:A:20:THR:HG23	1:A:30:ILE:HD13	1.80	0.64
3:C:1002:P6G:H62	4:C:1117:HOH:O	1.96	0.64
1:D:80:LYS:HZ2	3:D:1002:P6G:H91	1.66	0.60
1:A:59:ARG:HD3	1:B:57[A]:THR:CG2	2.31	0.60
1:C:29:PHE:C	1:C:29:PHE:CD1	2.78	0.60
1:C:80:LYS:NZ	3:C:1002:P6G:H91	2.15	0.60
1:C:95:GLN:NE2	1:D:114:THR:H	1.93	0.60
1:C:80:LYS:NZ	3:C:1002:P6G:H141	2.18	0.59
1:D:134:LYS:HA	1:D:134:LYS:NZ	2.16	0.58
1:D:80:LYS:NZ	3:D:1002:P6G:H141	2.19	0.58
1:A:57[B]:THR:HG21	4:B:1116:HOH:O	2.04	0.56
3:C:1002:P6G:H152	1:D:63:ALA:HB2	1.87	0.56
1:B:103:LYS:NZ	1:B:131:THR:HG22	2.22	0.54
1:A:59:ARG:HG2	1:B:57[B]:THR:HG22	1.90	0.53
1:D:115:THR:HG23	1:D:117:ALA:H	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57[B]:THR:HG23	1:B:59:ARG:HD3	1.92	0.51
1:D:80:LYS:HZ1	3:D:1002:P6G:H141	1.74	0.50
1:A:20:THR:HG23	1:A:30:ILE:HD11	1.91	0.50
1:C:90:THR:OG1	1:C:110:LEU:HD13	2.12	0.50
1:C:29:PHE:HE1	1:C:31:VAL:CG1	2.24	0.50
1:B:29[B]:PHE:CD2	1:B:130:PHE:HE2	2.30	0.49
1:A:91:THR:HB	1:B:91:THR:HB	1.94	0.49
1:C:80:LYS:HZ3	3:C:1002:P6G:H91	1.77	0.49
1:C:80:LYS:HZ2	3:C:1002:P6G:H141	1.76	0.48
1:D:80:LYS:NZ	3:D:1002:P6G:H122	2.28	0.48
1:C:63:ALA:CB	3:D:1002:P6G:H152	2.43	0.47
1:C:73:LEU:HD12	1:C:73:LEU:C	2.39	0.47
1:B:45:ASN:HB3	1:B:52:SER:HB2	1.97	0.46
1:A:21:TRP:HB2	1:A:29[A]:PHE:HD2	1.80	0.46
4:C:1123:HOH:O	1:D:57[A]:THR:HG21	2.16	0.46
1:C:57:THR:HG22	1:D:59:ARG:HG2	1.99	0.45
1:B:29[A]:PHE:HD2	1:B:56:LEU:CD2	2.29	0.45
1:D:115:THR:HG22	1:D:118:ASN:H	1.83	0.44
1:A:105:ASN:OD1	1:A:129[B]:THR:CG2	2.63	0.43
1:D:80:LYS:HZ2	3:D:1002:P6G:C9	2.30	0.43
1:B:21:TRP:HB2	1:B:29[A]:PHE:CD1	2.53	0.43
1:D:44:GLU:HB3	1:D:53:ARG:HG2	2.00	0.43
1:D:74:GLY:HA2	1:D:92:TRP:O	2.18	0.43
1:A:23:ASP:HB2	1:A:128:ASP:CG	2.44	0.43
1:B:21:TRP:HB2	1:B:29[A]:PHE:HD1	1.84	0.42
1:D:23:ASP:HB2	1:D:128:ASP:CG	2.44	0.42
1:C:81:ASN:HB2	3:C:1002:P6G:H22	2.00	0.42
1:B:23:ASP:HB3	1:B:130:PHE:CE2	2.55	0.41
1:D:80:LYS:HZ2	3:D:1002:P6G:C12	2.29	0.41
1:C:51:GLU:HB3	1:C:84:LYS:NZ	2.36	0.40
1:C:74:GLY:HA2	1:C:92:TRP:O	2.21	0.40
1:C:56:LEU:HD12	1:C:56:LEU:C	2.47	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	129/147 (88%)	124 (96%)	5 (4%)	0	100	100
1	B	128/147 (87%)	122 (95%)	6 (5%)	0	100	100
1	C	123/147 (84%)	119 (97%)	4 (3%)	0	100	100
1	D	124/147 (84%)	121 (98%)	3 (2%)	0	100	100
All	All	504/588 (86%)	486 (96%)	18 (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	97/107 (91%)	90 (93%)	7 (7%)	13	3
1	B	96/107 (90%)	87 (91%)	9 (9%)	8	1
1	C	91/107 (85%)	88 (97%)	3 (3%)	33	17
1	D	92/107 (86%)	87 (95%)	5 (5%)	20	6
All	All	376/428 (88%)	352 (94%)	24 (6%)	17	4

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	SER
1	A	44	GLU
1	A	47	VAL
1	A	49	ASN
1	A	51	GLU
1	A	52	SER
1	A	73	LEU
1	B	29[A]	PHE

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Mol	Chain	Res	Type
1	B	29[B]	PHE
1	B	44	GLU
1	B	47	VAL
1	B	57[A]	THR
1	B	57[B]	THR
1	B	73	LEU
1	B	115	THR
1	B	133	VAL
1	C	22	SER
1	C	29	PHE
1	C	44	GLU
1	D	12	SER
1	D	44	GLU
1	D	73	LEU
1	D	115	THR
1	D	134	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	A	95	GLN
1	A	107	GLN
1	A	116	ASN
1	B	49	ASN
1	B	116	ASN
1	C	95	GLN
1	C	107	GLN
1	D	116	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	D	1001	-	24,25,25	1.01	1 (4%)	28,32,32	1.68	5 (17%)
3	P6G	C	1002	-	15,15,18	0.96	0	14,14,17	1.12	1 (7%)
2	ZOF	B	1001	-	24,25,25	1.45	4 (16%)	28,32,32	1.57	4 (14%)
2	ZOF	A	1001	-	24,25,25	1.18	2 (8%)	28,32,32	1.77	5 (17%)
3	P6G	D	1002	-	15,15,18	0.81	0	14,14,17	1.03	2 (14%)
2	ZOF	C	1001	-	24,25,25	0.98	1 (4%)	28,32,32	1.52	5 (17%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ZOF	D	1001	-	-	4/16/37/37	0/2/2/2
3	P6G	C	1002	-	-	7/13/13/16	-
2	ZOF	B	1001	-	-	4/16/37/37	0/2/2/2
2	ZOF	A	1001	-	-	4/16/37/37	0/2/2/2
3	P6G	D	1002	-	-	6/13/13/16	-
2	ZOF	C	1001	-	-	5/16/37/37	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1001	ZOF	CAC-NAB	3.90	1.51	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1001	ZOF	CAF-CAC	3.37	1.60	1.53
2	A	1001	ZOF	CAF-CAC	2.69	1.58	1.53
2	A	1001	ZOF	CAI-CAH	2.59	1.59	1.52
2	B	1001	ZOF	CAU-CAV	2.26	1.55	1.50
2	C	1001	ZOF	CAU-CAV	2.21	1.55	1.50
2	D	1001	ZOF	CAF-CAC	2.15	1.57	1.53
2	B	1001	ZOF	CAI-CAH	2.06	1.58	1.52

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1001	ZOF	NAE-CAA-NAB	5.33	114.63	109.21
2	A	1001	ZOF	NAE-CAA-NAB	5.19	114.49	109.21
2	B	1001	ZOF	NAE-CAA-NAB	4.76	114.05	109.21
2	C	1001	ZOF	NAE-CAA-NAB	4.28	113.56	109.21
2	D	1001	ZOF	CAC-CAF-SAG	-3.72	100.94	106.06
2	A	1001	ZOF	CAK-CAI-CAH	-3.52	105.94	114.04
2	A	1001	ZOF	CAH-CAD-NAE	-3.35	109.80	113.34
2	B	1001	ZOF	CAK-CAI-CAH	-3.19	106.70	114.04
2	C	1001	ZOF	CAC-CAF-SAG	-3.17	101.70	106.06
2	C	1001	ZOF	CAK-CAI-CAH	-2.86	107.46	114.04
2	D	1001	ZOF	CAK-CAI-CAH	-2.73	107.75	114.04
2	B	1001	ZOF	CAC-CAF-SAG	-2.35	102.82	106.06
2	B	1001	ZOF	CAC-CAD-NAE	2.27	105.24	102.68
3	D	1002	P6G	O13-C14-C15	2.21	119.86	110.11
2	C	1001	ZOF	CAH-CAD-NAE	-2.19	111.02	113.34
2	D	1001	ZOF	OAX-CAV-CAU	-2.14	116.30	123.09
2	A	1001	ZOF	OAX-CAV-CAU	-2.14	116.31	123.09
2	A	1001	ZOF	CAT-CAU-CAV	-2.09	109.05	114.51
2	C	1001	ZOF	CAC-CAD-NAE	2.05	104.99	102.68
3	D	1002	P6G	O4-C3-C2	2.04	119.10	110.11
2	D	1001	ZOF	CAC-CAD-NAE	2.03	104.97	102.68
3	C	1002	P6G	O4-C3-C2	2.01	118.95	110.11

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	1001	ZOF	CAR-CAQ-NAO-CAN
3	C	1002	P6G	O7-C8-C9-O10
3	D	1002	P6G	O7-C8-C9-O10
3	C	1002	P6G	O4-C5-C6-O7

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Mol	Chain	Res	Type	Atoms
2	C	1001	ZOF	CAS-CAT-CAU-CAV
3	D	1002	P6G	O4-C5-C6-O7
3	D	1002	P6G	O1-C2-C3-O4
2	B	1001	ZOF	CAS-CAT-CAU-CAV
3	C	1002	P6G	O1-C2-C3-O4
2	D	1001	ZOF	CAS-CAT-CAU-CAV
2	A	1001	ZOF	CAS-CAT-CAU-CAV
2	A	1001	ZOF	CAQ-CAR-CAS-CAT
3	C	1002	P6G	O13-C14-C15-O16
3	D	1002	P6G	O13-C14-C15-O16
3	C	1002	P6G	O10-C11-C12-O13
2	B	1001	ZOF	CAQ-CAR-CAS-CAT
3	D	1002	P6G	O10-C11-C12-O13
3	D	1002	P6G	C15-C14-O13-C12
3	C	1002	P6G	C15-C14-O13-C12
2	D	1001	ZOF	CAT-CAU-CAV-OAX
2	C	1001	ZOF	CAT-CAU-CAV-OAW
2	B	1001	ZOF	CAT-CAU-CAV-OAW
2	A	1001	ZOF	CAT-CAU-CAV-OAW
2	D	1001	ZOF	CAT-CAU-CAV-OAW
2	B	1001	ZOF	CAT-CAU-CAV-OAX
2	A	1001	ZOF	CAT-CAU-CAV-OAX
2	C	1001	ZOF	CAT-CAU-CAV-OAX
2	C	1001	ZOF	SAG-CAH-CAI-CAK
2	C	1001	ZOF	CAR-CAQ-NAO-CAN
3	C	1002	P6G	C11-C12-O13-C14

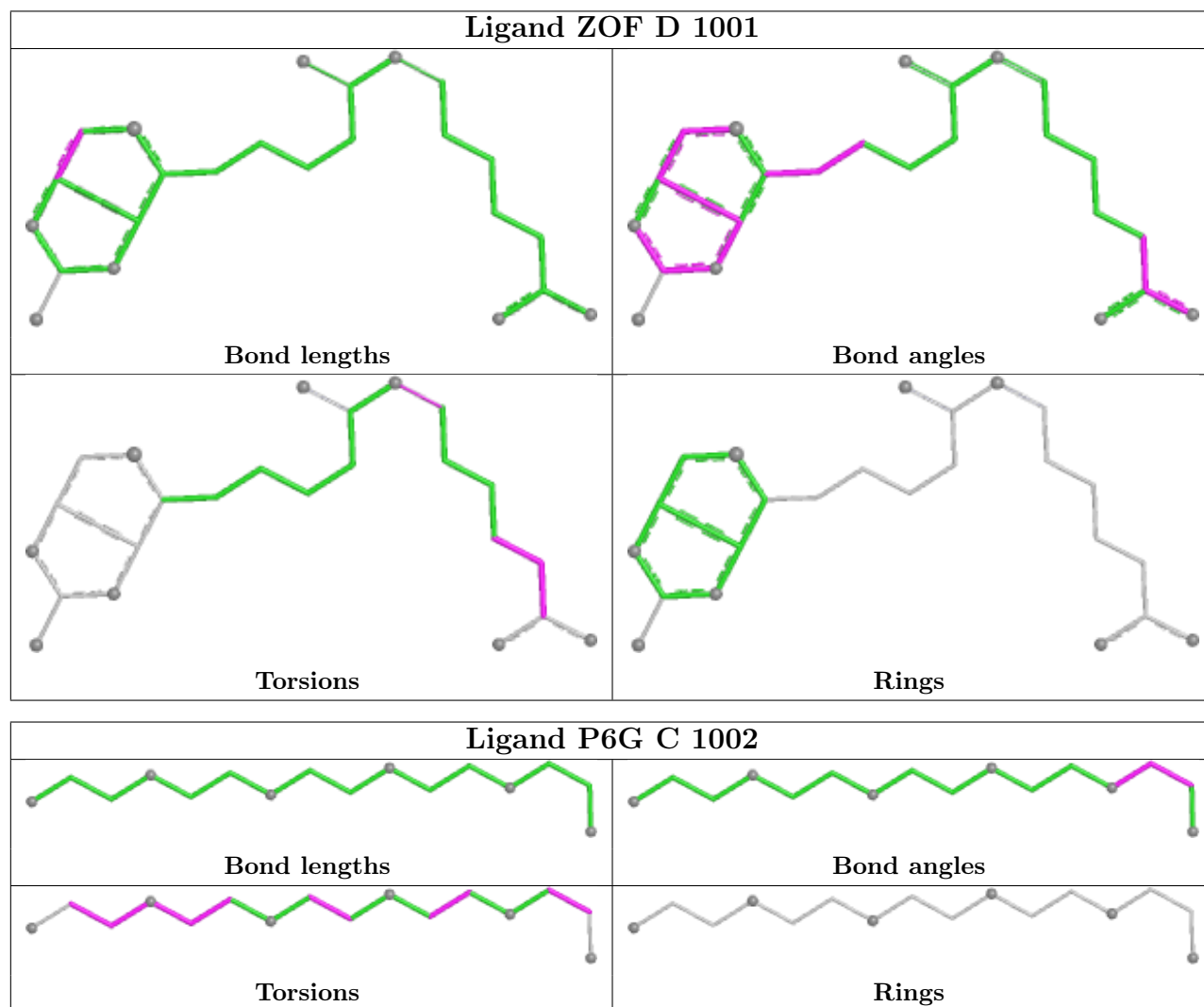
There are no ring outliers.

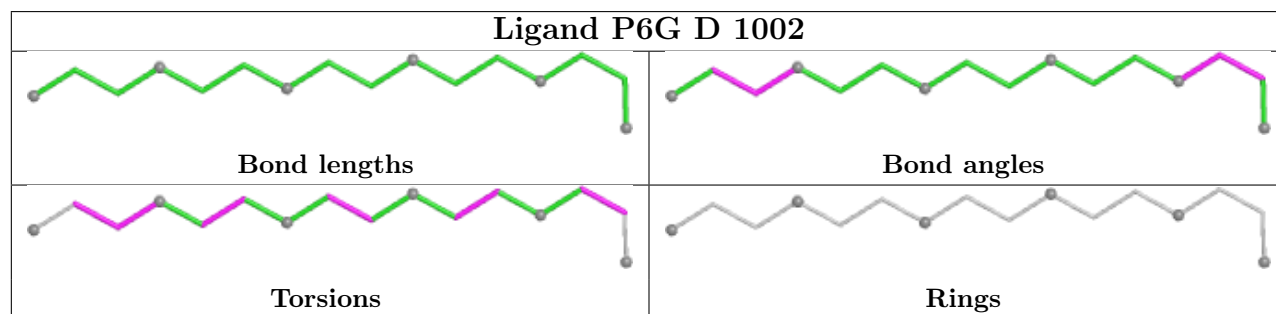
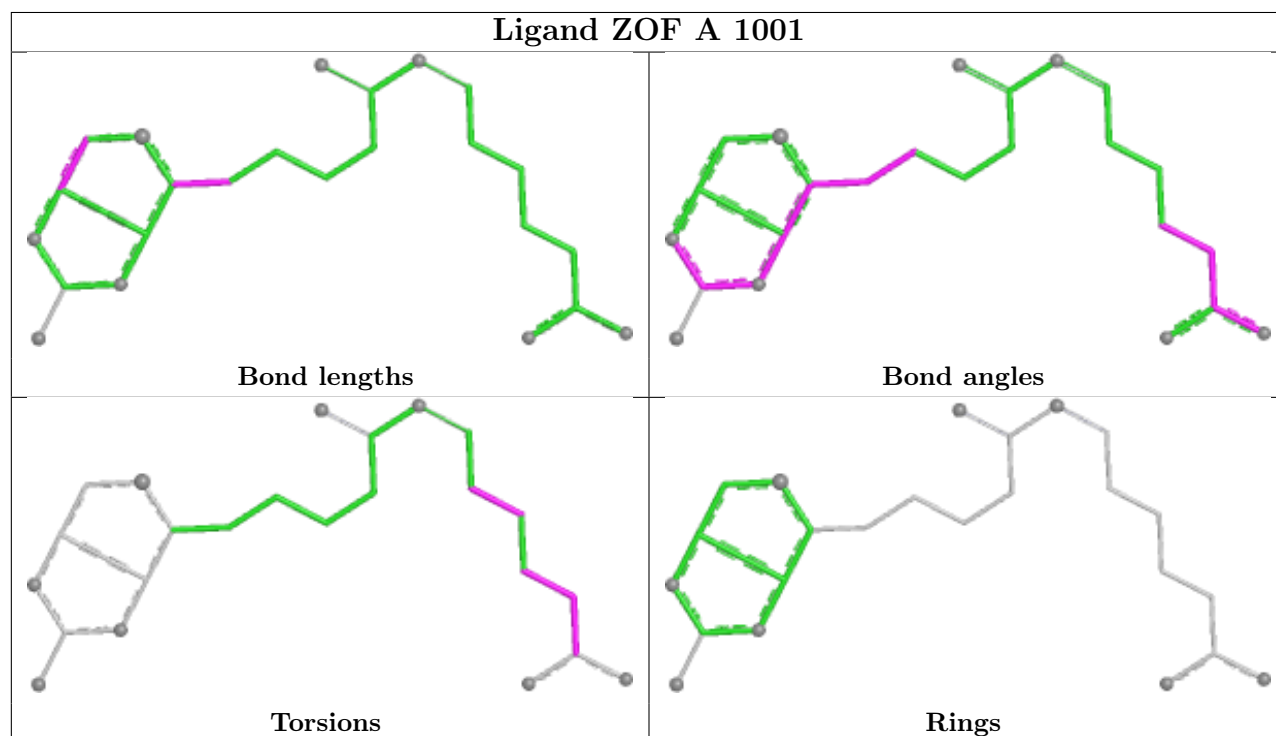
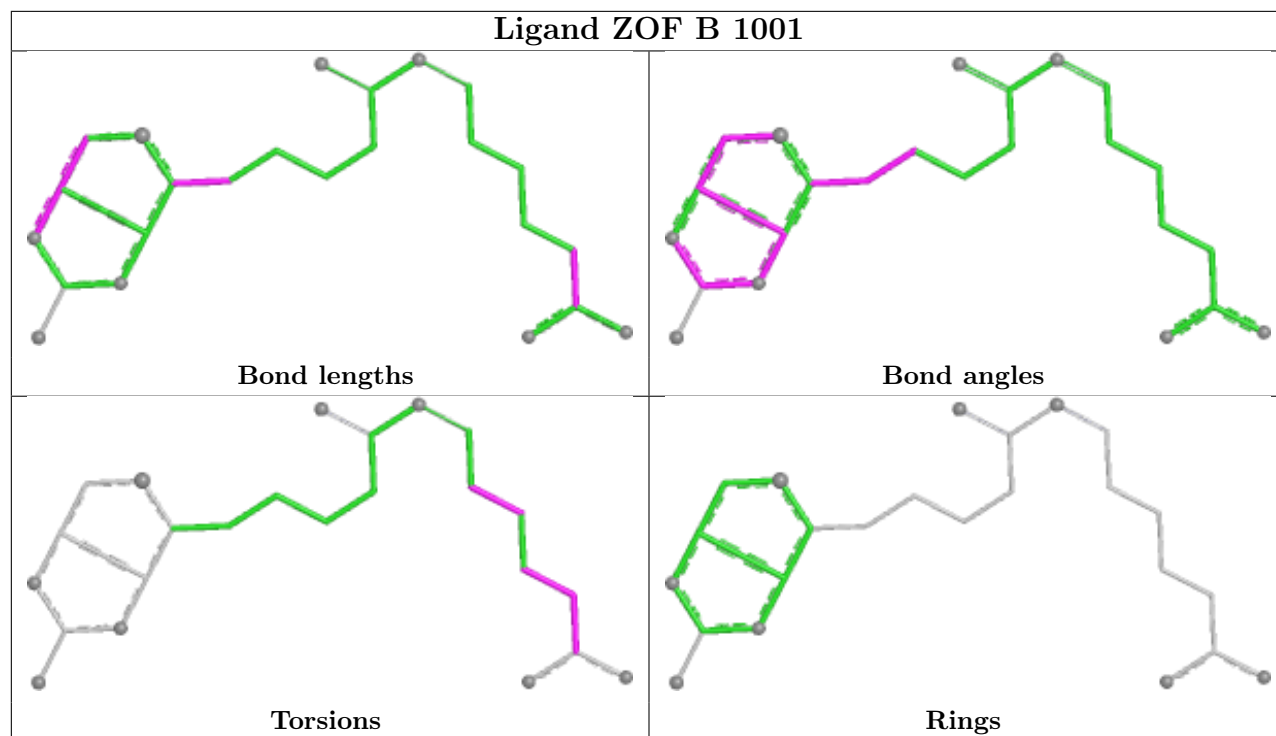
2 monomers are involved in 17 short contacts:

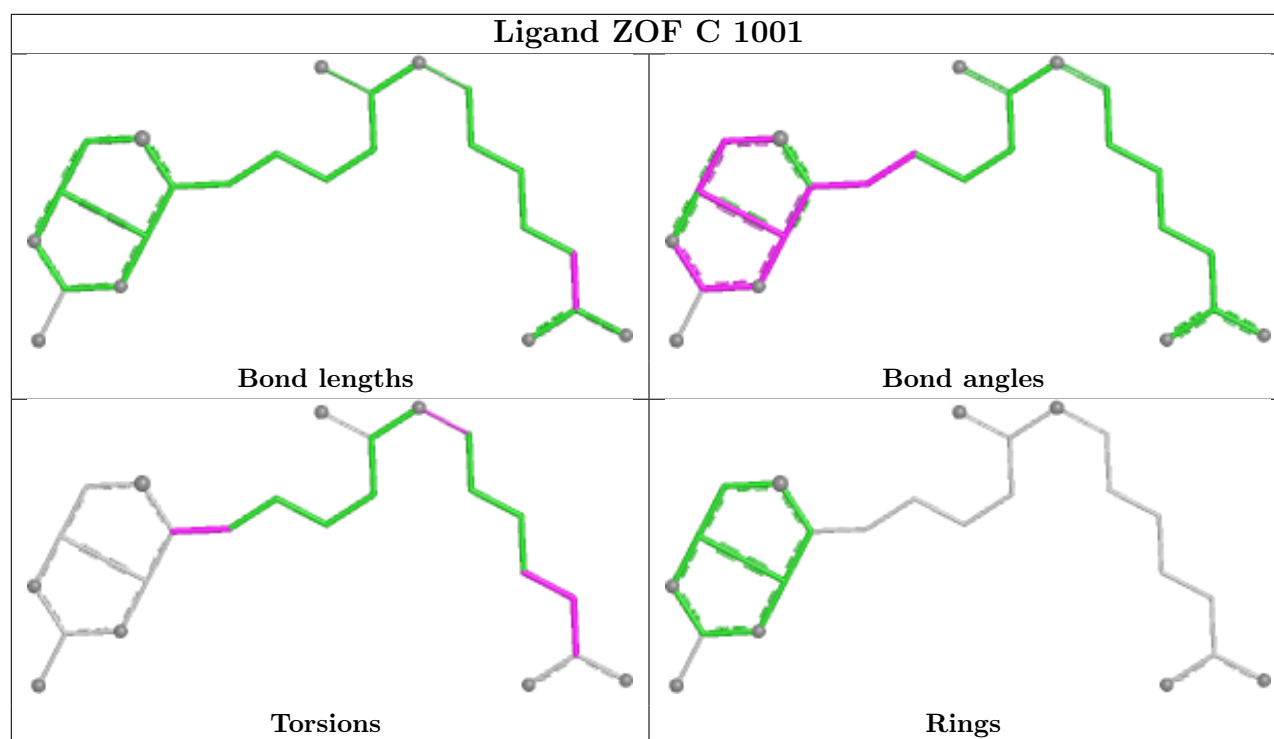
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1002	P6G	7	0
3	D	1002	P6G	10	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	125/147 (85%)	-1.05	0 100 100	12, 23, 49, 63	6 (4%)
1	B	125/147 (85%)	-1.10	0 100 100	12, 22, 50, 63	5 (4%)
1	C	124/147 (84%)	-1.04	0 100 100	14, 25, 47, 55	1 (0%)
1	D	125/147 (85%)	-1.03	0 100 100	12, 26, 47, 58	1 (0%)
All	All	499/588 (84%)	-1.06	0 100 100	12, 24, 49, 63	13 (2%)

There are no RSRZ outliers to report.

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.

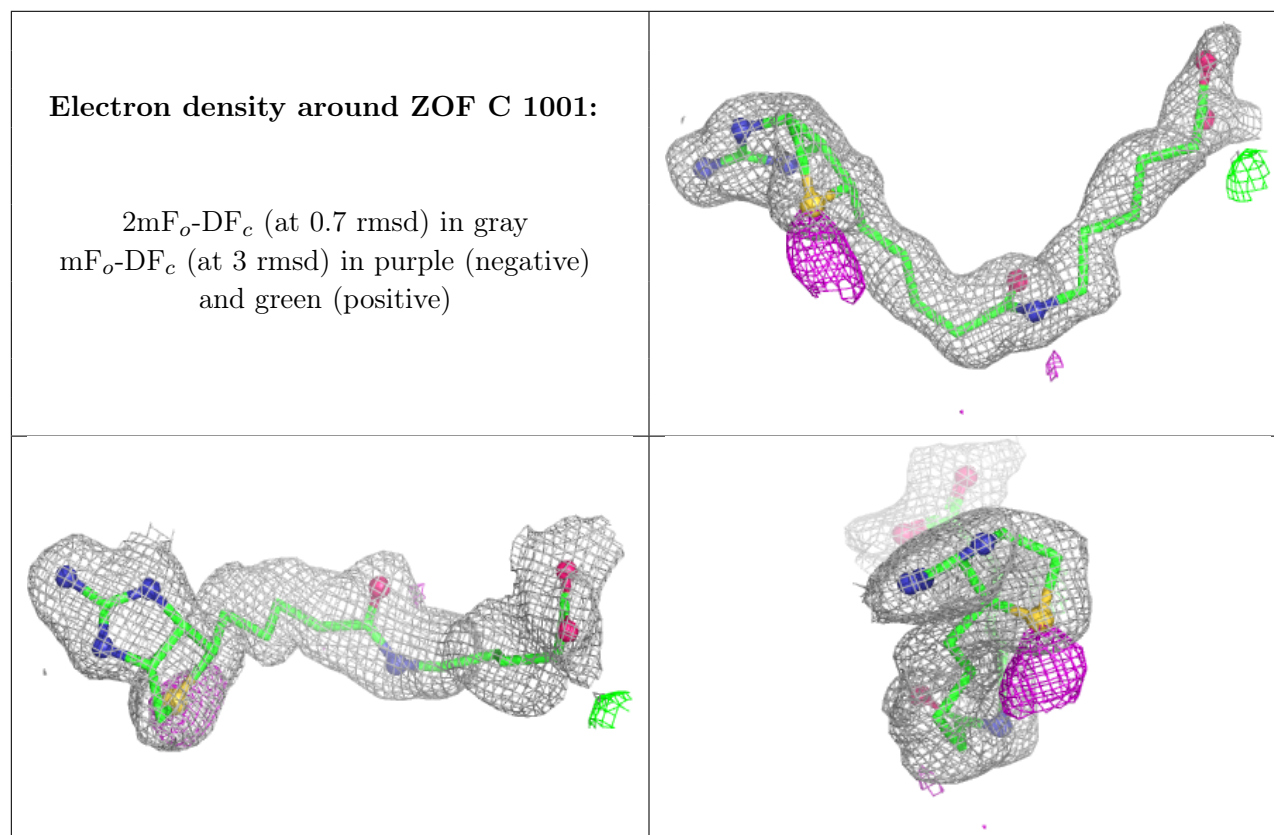
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ZOF	C	1001	24/24	0.98	0.04	26,33,61,64	0
3	P6G	C	1002	16/19	0.98	0.05	42,45,50,50	0
3	P6G	D	1002	16/19	0.98	0.04	35,46,51,51	0
2	ZOF	D	1001	24/24	0.99	0.04	26,35,63,65	0

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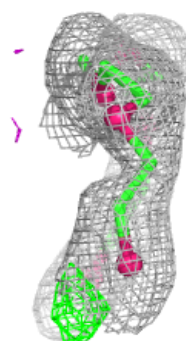
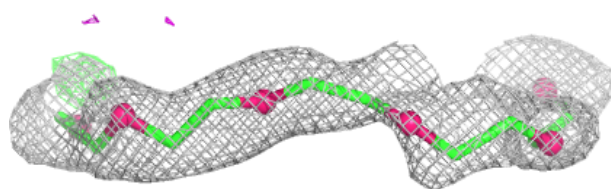
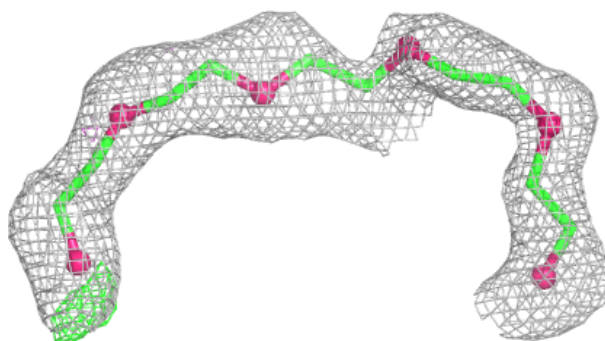
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZOF	B	1001	24/24	0.99	0.04	21,31,56,57	0
2	ZOF	A	1001	24/24	0.99	0.04	23,29,63,65	0

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.

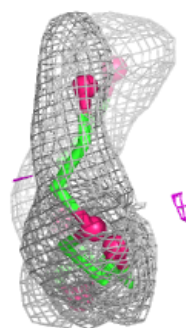
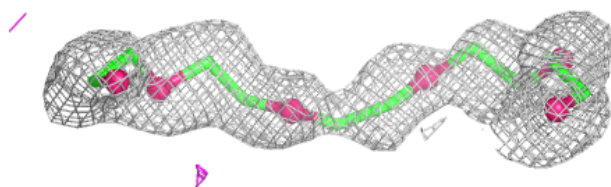
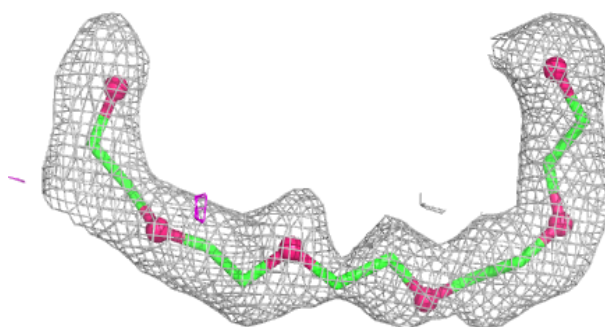


Electron density around P6G C 1002:

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

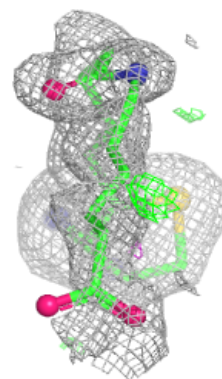
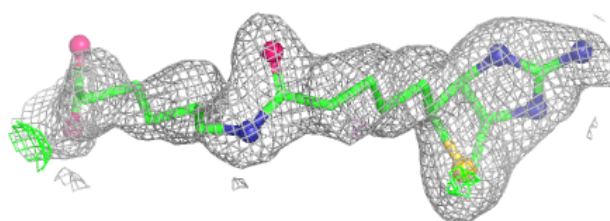
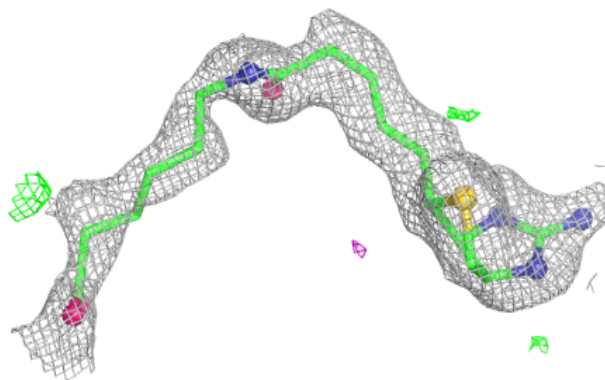
**Electron density around P6G D 1002:**

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

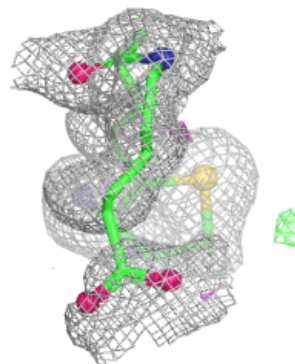
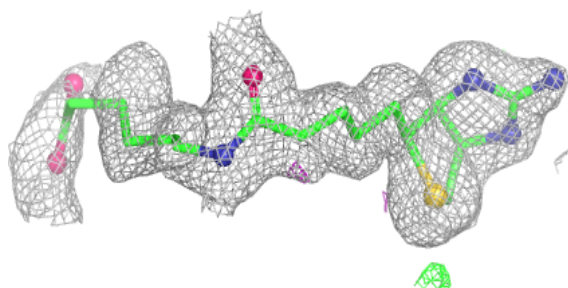
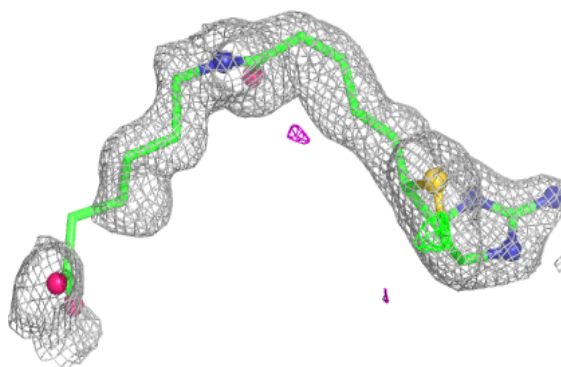


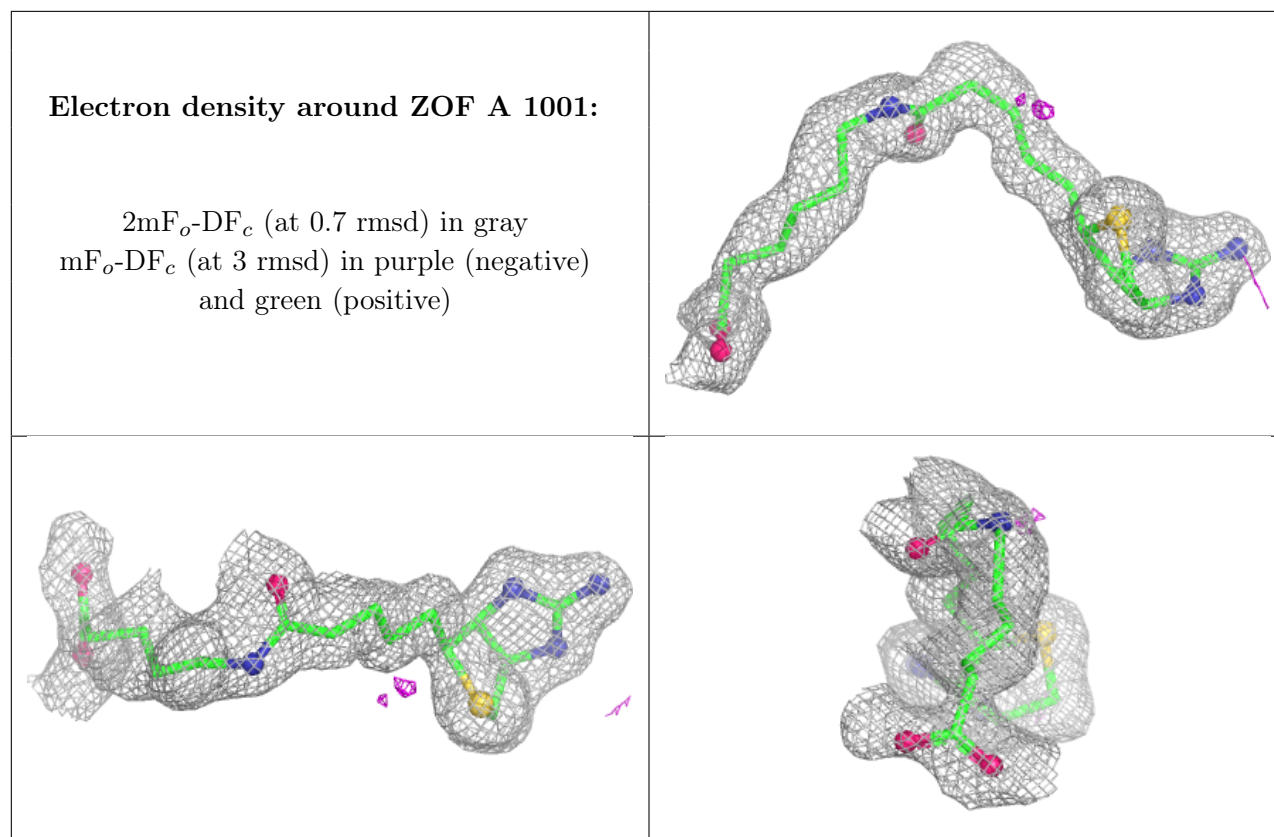
Electron density around ZOF D 1001:

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

**Electron density around ZOF B 1001:**

$2mF_o-DF_c$ (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.