



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 07:08 AM UTC

PDB ID : 5Z97 / pdb_00005z97
Title : Crystal structure of a lactonase double mutant in complex with ligand N
Authors : Zheng, Y.Y.; Liu, W.D.; Chen, C.C.; Guo, R.T.
Deposited on : 2018-02-02
Resolution : 2.32 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

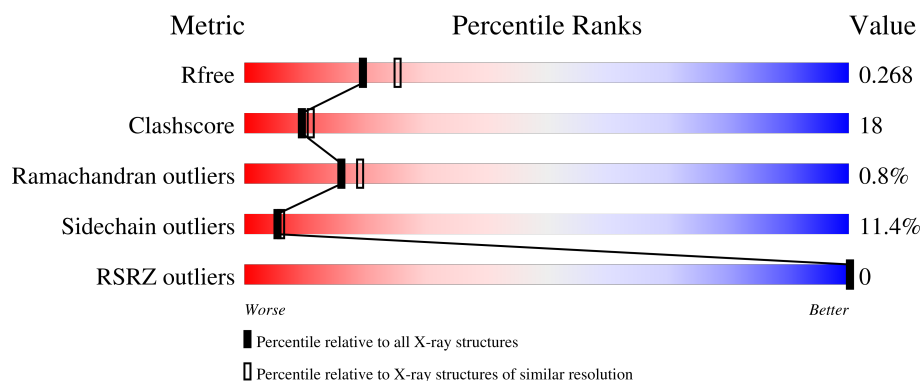
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.32 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	7754 (2.34-2.30)
Clashscore	190562	8383 (2.34-2.30)
Ramachandran outliers	187476	8303 (2.34-2.30)
Sidechain outliers	187428	8303 (2.34-2.30)
RSRZ outliers	180081	7760 (2.34-2.30)

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	263	 77% 21% .
1	B	263	 79% 20% .
1	C	263	 71% 23% 6%
1	D	263	 40% 44% 14% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8498 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.

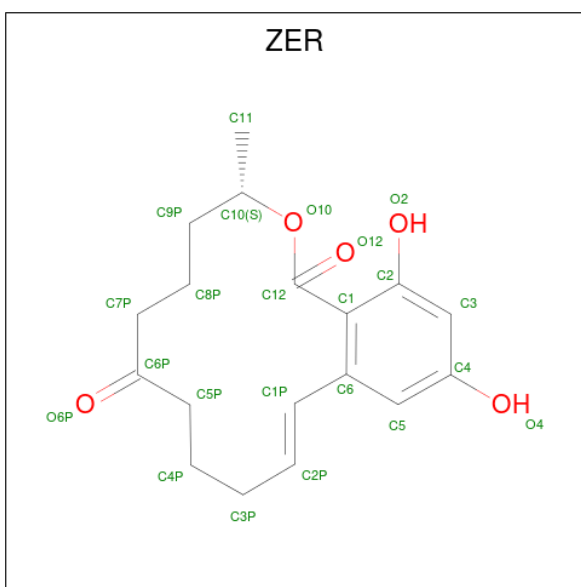
- Molecule 1 is a protein called Lactonase for protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			2048	1300	347	392	9			
1	B	263	Total	C	N	O	S	0	0	0
			2048	1300	347	392	9			
1	C	263	Total	C	N	O	S	0	0	0
			2048	1300	347	392	9			
1	D	261	Total	C	N	O	S	0	0	0
			2030	1288	344	389	9			

There are 8 discrepancies between the modelled and reference sequences:

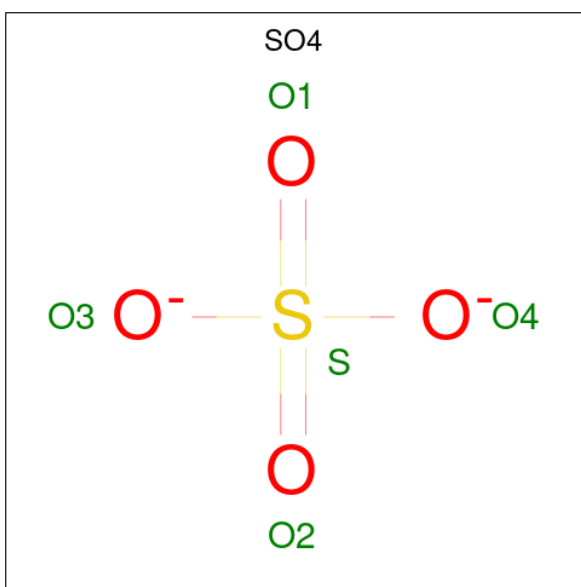
Chain	Residue	Modelled	Actual	Comment	Reference
A	105	ALA	SER	engineered mutation	UNP A0A0D2ILK1
A	160	ALA	TYR	engineered mutation	UNP A0A0D2ILK1
B	105	ALA	SER	engineered mutation	UNP A0A0D2ILK1
B	160	ALA	TYR	engineered mutation	UNP A0A0D2ILK1
C	105	ALA	SER	engineered mutation	UNP A0A0D2ILK1
C	160	ALA	TYR	engineered mutation	UNP A0A0D2ILK1
D	105	ALA	SER	engineered mutation	UNP A0A0D2ILK1
D	160	ALA	TYR	engineered mutation	UNP A0A0D2ILK1

- Molecule 2 is (3S,11E)-14,16-dihydroxy-3-methyl-3,4,5,6,9,10-hexahydro-1H-2-benzoxacyclo tetradecine-1,7(8H)-dione (CCD ID: ZER) (formula: C₁₈H₂₂O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			23	18	5		
2	B	1	Total	C	O	0	0
			23	18	5		
2	C	1	Total	C	O	0	0
			23	18	5		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		

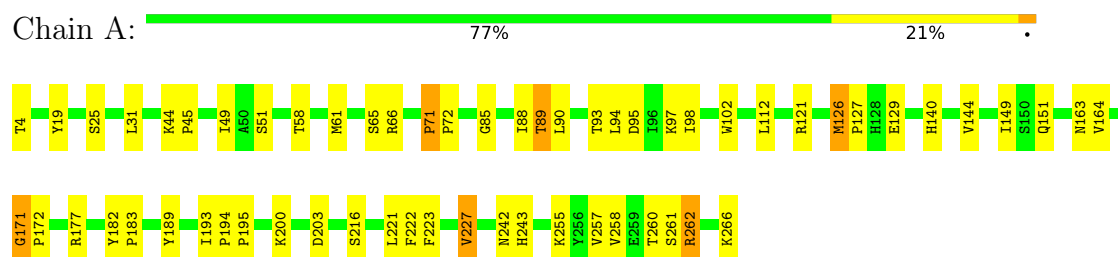
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	74	Total 74	O 74	0	0
4	B	97	Total 97	O 97	0	0
4	C	59	Total 59	O 59	0	0
4	D	20	Total 20	O 20	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.

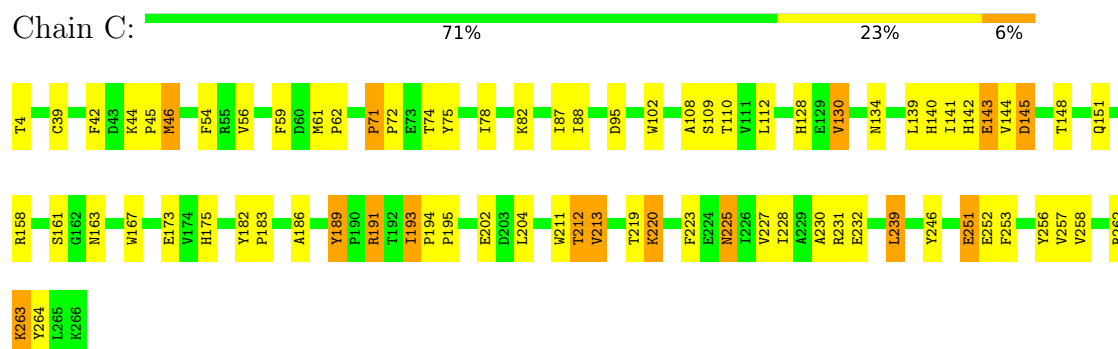
- Molecule 1: Lactonase for protein



- Molecule 1: Lactonase for protein

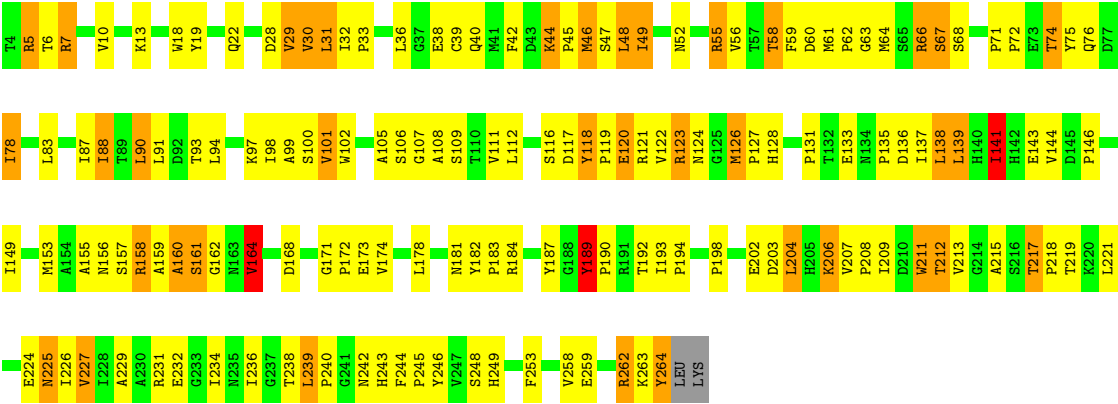


- Molecule 1: Lactonase for protein



- Molecule 1: Lactonase for protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	44.32Å 110.52Å 127.78Å 90.00° 99.93° 90.00°	Depositor
Resolution (Å)	25.00 – 2.32 25.00 – 2.32	Depositor EDS
% Data completeness (in resolution range)	94.9 (25.00-2.32) 94.7 (25.00-2.32)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.210 , 0.266 0.217 , 0.268	Depositor DCC
R_{free} test set	2490 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	40.1	Xtriage
Anisotropy	0.022	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 29.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.277 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8498	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.25% 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: SO4, ZER

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	0.76	1/2106 (0.0%)	0.93	7/2875 (0.2%)
1	B	0.72	0/2106	0.87	4/2875 (0.1%)
1	C	0.58	0/2106	0.94	9/2875 (0.3%)
1	D	0.62	0/2088	1.06	9/2853 (0.3%)
All	All	0.67	1/8406 (0.0%)	0.95	29/11478 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	164	VAL	C-O	-5.25	1.17	1.24

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	171	GLY	CA-C-N	11.12	132.00	119.32
1	D	171	GLY	C-N-CA	11.12	132.00	119.32
1	C	145	ASP	CA-C-N	8.45	128.03	119.24
1	C	145	ASP	C-N-CA	8.45	128.03	119.24
1	D	161	SER	N-CA-C	8.34	120.45	111.36
1	A	171	GLY	CA-C-N	7.93	127.85	119.05
1	A	171	GLY	C-N-CA	7.93	127.85	119.05
1	C	71	PRO	CA-C-N	7.67	128.06	119.32
1	C	71	PRO	C-N-CA	7.67	128.06	119.32
1	D	126	MET	CA-C-N	6.93	127.48	120.14
1	D	126	MET	C-N-CA	6.93	127.48	120.14
1	C	130	VAL	CB-CA-C	-6.91	104.52	110.94
1	D	141	ILE	N-CA-C	6.90	121.53	112.76
1	A	71	PRO	CA-C-N	6.50	127.01	119.47
1	A	71	PRO	C-N-CA	6.50	127.01	119.47
1	D	160	ALA	N-CA-C	5.95	119.50	112.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	65	SER	CB-CA-C	-5.84	109.82	116.54
1	D	189	TYR	CA-C-N	5.83	127.13	119.84
1	D	189	TYR	C-N-CA	5.83	127.13	119.84
1	C	239	LEU	CA-C-N	-5.66	114.42	120.03
1	C	239	LEU	C-N-CA	-5.66	114.42	120.03
1	B	71	PRO	CA-C-N	5.63	126.88	119.84
1	B	71	PRO	C-N-CA	5.63	126.88	119.84
1	A	61	MET	CA-C-N	5.42	125.72	119.93
1	A	61	MET	C-N-CA	5.42	125.72	119.93
1	B	145	ASP	CA-C-N	5.37	124.55	118.97
1	B	145	ASP	C-N-CA	5.37	124.55	118.97
1	C	193	ILE	CA-C-N	5.00	125.53	120.38
1	C	193	ILE	C-N-CA	5.00	125.53	120.38

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	2048	0	1989	34	0
1	B	2048	0	1989	33	0
1	C	2048	0	1989	56	0
1	D	2030	0	1965	170	0
2	A	23	0	21	3	0
2	B	23	0	21	0	0
2	C	23	0	21	3	0
3	A	5	0	0	0	0
4	A	74	0	0	2	0
4	B	97	0	0	2	0
4	C	59	0	0	2	0
4	D	20	0	0	2	0
All	All	8498	0	7995	290	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 18.

All (290) 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:75:TYR:C	1:D:78:ILE:HD11	1.76	1.10
1:D:135:PRO:HD2	1:D:138:LEU:HD11	1.23	1.09
1:D:162:GLY:HA3	1:D:248:SER:HB2	1.44	0.96
1:D:49:ILE:CD1	1:D:258:VAL:HG13	1.95	0.94
1:D:160:ALA:O	1:D:242:ASN:HB2	1.66	0.94
1:A:126:MET:SD	1:A:260:THR:HG21	2.07	0.93
1:C:212:THR:HG21	1:C:239:LEU:HG	1.51	0.93
1:D:99:ALA:O	1:D:123:ARG:HB3	1.67	0.93
1:C:230:ALA:HB2	1:D:227:VAL:HG12	1.51	0.90
1:D:49:ILE:HD13	1:D:258:VAL:HG13	1.53	0.90
1:D:62:PRO:HB2	1:D:74:THR:CG2	2.04	0.88
1:D:135:PRO:HD2	1:D:138:LEU:CD1	2.02	0.88
1:D:124:ASN:ND2	1:D:264:TYR:CE1	2.42	0.88
1:D:207:VAL:HB	1:D:208:PRO:HD2	1.53	0.87
1:D:62:PRO:HB2	1:D:74:THR:HG23	1.58	0.86
1:D:62:PRO:CB	1:D:74:THR:CG2	2.53	0.85
1:C:223:PHE:O	1:C:227:VAL:HG23	1.77	0.84
1:D:66:ARG:HG2	1:D:66:ARG:HH21	1.42	0.84
1:D:135:PRO:O	1:D:139:LEU:HD12	1.77	0.84
1:D:75:TYR:CA	1:D:78:ILE:HD11	2.07	0.84
1:D:227:VAL:O	1:D:231:ARG:HG3	1.78	0.83
1:B:252:GLU:N	1:B:252:GLU:OE1	2.12	0.83
1:D:28:ASP:HA	1:D:55:ARG:O	1.80	0.82
1:D:29:VAL:HG22	1:D:100:SER:HB2	1.61	0.82
1:D:99:ALA:C	1:D:123:ARG:HB3	2.04	0.82
1:D:62:PRO:HB3	1:D:74:THR:HG21	1.62	0.82
1:D:29:VAL:O	1:D:56:VAL:HA	1.79	0.81
1:D:28:ASP:OD1	1:D:55:ARG:HG2	1.80	0.81
1:D:49:ILE:CD1	1:D:258:VAL:CG1	2.59	0.80
1:D:207:VAL:HB	1:D:208:PRO:CD	2.11	0.79
1:D:135:PRO:HG3	1:D:221:LEU:HD21	1.65	0.79
1:D:39:CYS:HB2	1:D:58:THR:HG23	1.62	0.79
1:D:72:PRO:HA	1:D:75:TYR:CZ	2.18	0.78
1:C:252:GLU:N	1:C:252:GLU:OE1	2.17	0.77
1:A:243:HIS:CD2	2:A:301:ZER:H6	2.20	0.77
1:D:137:ILE:HG13	1:D:138:LEU:HG	1.67	0.77
1:D:160:ALA:O	1:D:243:HIS:N	2.13	0.77
1:D:75:TYR:HA	1:D:78:ILE:HD11	1.66	0.76
1:D:158:ARG:HD2	1:D:158:ARG:C	2.10	0.76
1:C:223:PHE:CE1	1:C:227:VAL:HG21	2.22	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:29:VAL:O	1:D:56:VAL:HG23	1.86	0.75
1:C:142:HIS:CD2	1:C:191:ARG:HD3	2.23	0.74
1:D:66:ARG:HG2	1:D:66:ARG:NH2	1.98	0.74
1:D:135:PRO:HB2	1:D:137:ILE:HG12	1.70	0.74
1:D:141:ILE:HD13	1:D:149:ILE:HG12	1.68	0.74
1:B:260:THR:O	1:B:263:LYS:HG2	1.87	0.73
1:D:39:CYS:O	1:D:42:PHE:HB2	1.89	0.73
1:D:88:ILE:HG13	1:D:118:TYR:CD2	2.24	0.73
1:C:42:PHE:O	1:C:46:MET:HB2	1.90	0.72
1:D:28:ASP:OD1	1:D:55:ARG:HB3	1.90	0.72
1:D:39:CYS:HB2	1:D:58:THR:CG2	2.18	0.72
1:D:182:TYR:HB2	1:D:183:PRO:HD3	1.70	0.72
1:D:107:GLY:O	1:D:111:VAL:HG23	1.89	0.72
1:D:62:PRO:HB3	1:D:74:THR:CG2	2.20	0.71
1:D:75:TYR:C	1:D:78:ILE:CD1	2.61	0.71
1:D:211:TRP:HB3	1:D:236:ILE:HD12	1.72	0.71
1:D:30:VAL:C	1:D:31:LEU:HD23	2.16	0.70
1:D:112:LEU:HB2	1:D:198:PRO:HG2	1.74	0.70
1:D:135:PRO:O	1:D:139:LEU:CD1	2.38	0.70
1:D:29:VAL:N	1:D:55:ARG:O	2.24	0.70
1:D:135:PRO:CD	1:D:138:LEU:HD11	2.14	0.69
1:D:36:LEU:O	1:D:181:ASN:ND2	2.25	0.69
1:D:44:LYS:HB2	1:D:45:PRO:CD	2.23	0.69
1:D:60:ASP:OD2	1:D:66:ARG:HB2	1.93	0.69
1:D:62:PRO:CB	1:D:74:THR:HG23	2.19	0.69
1:D:100:SER:OG	1:D:123:ARG:HG2	1.93	0.69
1:D:88:ILE:HD13	1:D:91:LEU:HD12	1.74	0.68
1:D:62:PRO:CB	1:D:74:THR:HG21	2.23	0.67
1:D:120:GLU:CD	1:D:120:GLU:H	2.02	0.67
1:D:217:THR:HG22	1:D:218:PRO:HD2	1.76	0.67
1:B:144:VAL:HG23	1:B:149:ILE:HG13	1.77	0.66
1:A:4:THR:HG22	1:C:151:GLN:OE1	1.94	0.66
1:D:10:VAL:HG21	1:D:94:LEU:HD21	1.77	0.66
1:D:52:ASN:OD1	1:D:262:ARG:NH2	2.29	0.66
1:D:7:ARG:HD2	1:D:19:TYR:OH	1.96	0.66
1:D:39:CYS:CB	1:D:58:THR:HG23	2.25	0.66
1:D:76:GLN:N	1:D:78:ILE:HD11	2.10	0.66
1:D:160:ALA:C	1:D:242:ASN:HB2	2.21	0.66
1:D:161:SER:O	1:D:248:SER:HB3	1.96	0.65
1:D:44:LYS:HB2	1:D:45:PRO:HD3	1.79	0.65
1:D:227:VAL:O	1:D:231:ARG:CG	2.44	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:MET:SD	1:A:260:THR:CG2	2.84	0.64
1:D:189:TYR:H	1:D:190:PRO:HD2	1.61	0.64
1:D:259:GLU:HA	1:D:262:ARG:HG3	1.79	0.64
1:D:28:ASP:OD1	1:D:55:ARG:CG	2.47	0.63
1:A:144:VAL:HG23	1:A:149:ILE:HG13	1.81	0.63
1:D:144:VAL:HG23	1:D:149:ILE:HG13	1.81	0.63
1:C:72:PRO:HA	1:C:75:TYR:CZ	2.35	0.61
1:C:220:LYS:HG2	1:D:215:ALA:HB1	1.83	0.61
1:D:207:VAL:CB	1:D:208:PRO:CD	2.77	0.61
1:A:257:VAL:O	1:A:260:THR:HG22	2.00	0.61
1:D:204:LEU:CD2	1:D:234:ILE:HD11	2.31	0.61
1:D:38:GLU:CD	1:D:40:GLN:HB3	2.25	0.60
1:D:206:LYS:HA	1:D:206:LYS:NZ	2.16	0.60
1:D:49:ILE:HD11	1:D:258:VAL:CG1	2.30	0.60
1:A:85:GLY:O	1:A:89:THR:HG23	2.01	0.60
1:D:39:CYS:HA	1:D:42:PHE:HD2	1.67	0.60
1:D:207:VAL:CB	1:D:208:PRO:HD2	2.28	0.60
2:C:301:ZER:O10	2:C:301:ZER:H11	2.01	0.59
1:D:49:ILE:HD12	1:D:49:ILE:O	2.02	0.59
1:D:155:ALA:O	1:D:158:ARG:HG3	2.03	0.59
1:A:4:THR:N	4:A:401:HOH:O	2.34	0.59
1:D:40:GLN:HG3	1:D:40:GLN:O	2.02	0.59
1:D:128:HIS:ND1	1:D:212:THR:O	2.35	0.59
1:D:66:ARG:HH21	1:D:66:ARG:CG	2.13	0.59
1:D:164:VAL:O	1:D:168:ASP:OD1	2.21	0.58
1:C:175:HIS:ND1	4:C:401:HOH:O	2.32	0.58
1:D:160:ALA:O	1:D:242:ASN:CB	2.48	0.58
1:A:243:HIS:HD2	2:A:301:ZER:H6	1.67	0.58
1:D:90:LEU:O	1:D:90:LEU:HG	1.99	0.58
1:A:260:THR:HG23	1:A:261:SER:N	2.18	0.58
1:D:204:LEU:HG	1:D:234:ILE:HD11	1.84	0.57
1:C:75:TYR:HA	1:C:78:ILE:HD11	1.87	0.57
1:D:39:CYS:SG	1:D:58:THR:HG23	2.44	0.57
1:D:116:SER:HB2	1:D:203:ASP:HB3	1.85	0.57
1:D:99:ALA:O	1:D:123:ARG:CB	2.48	0.57
1:D:88:ILE:HG13	1:D:118:TYR:CE2	2.40	0.56
1:D:28:ASP:OD1	1:D:55:ARG:CB	2.53	0.56
1:D:64:MET:HB3	1:D:181:ASN:ND2	2.20	0.56
1:B:128:HIS:CD2	1:B:129:GLU:HG3	2.40	0.56
1:D:52:ASN:OD1	1:D:52:ASN:O	2.23	0.56
1:D:153:MET:HE3	1:D:153:MET:HA	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:31:LEU:HD23	1:D:31:LEU:N	2.20	0.55
1:A:182:TYR:HB2	1:A:183:PRO:HD3	1.87	0.55
1:C:128:HIS:ND1	1:C:212:THR:HB	2.21	0.55
1:D:117:ASP:C	1:D:119:PRO:HD3	2.32	0.55
1:C:213:VAL:HG12	1:D:219:THR:HG21	1.87	0.54
1:C:102:TRP:CD1	1:C:102:TRP:C	2.85	0.54
1:D:63:GLY:HA3	1:D:75:TYR:CD2	2.42	0.54
1:B:231:ARG:NE	4:B:405:HOH:O	2.39	0.54
1:B:44:LYS:N	1:B:45:PRO:CD	2.70	0.54
1:C:182:TYR:HB2	1:C:183:PRO:HD3	1.90	0.54
1:C:134:ASN:ND2	2:C:301:ZER:O4	2.39	0.53
1:B:95:ASP:CG	1:B:95:ASP:O	2.51	0.53
1:D:161:SER:O	1:D:248:SER:CB	2.55	0.53
1:D:105:ALA:O	1:D:108:ALA:HB3	2.08	0.53
1:D:45:PRO:HA	1:D:48:LEU:HD22	1.91	0.53
1:D:238:THR:O	1:D:239:LEU:HD12	2.09	0.52
1:D:60:ASP:HB2	1:D:67:SER:OG	2.10	0.52
1:D:189:TYR:N	1:D:190:PRO:HD2	2.23	0.52
1:A:31:LEU:O	1:A:58:THR:OG1	2.24	0.52
1:D:136:ASP:HA	1:D:139:LEU:HD13	1.91	0.52
1:D:19:TYR:CD1	1:D:19:TYR:C	2.88	0.52
1:C:213:VAL:CG1	1:D:219:THR:HG21	2.41	0.51
1:D:206:LYS:HA	1:D:206:LYS:CE	2.40	0.51
1:D:118:TYR:N	1:D:118:TYR:CD1	2.79	0.51
1:C:142:HIS:CE1	1:C:143:GLU:HG2	2.47	0.50
1:D:215:ALA:N	1:D:239:LEU:O	2.45	0.50
1:C:158:ARG:HG3	1:C:167:TRP:CD1	2.46	0.50
1:D:71:PRO:O	1:D:74:THR:HB	2.12	0.50
1:C:72:PRO:HA	1:C:75:TYR:CE1	2.47	0.50
1:D:76:GLN:N	1:D:78:ILE:CD1	2.75	0.49
1:D:217:THR:CG2	1:D:218:PRO:HD2	2.41	0.49
1:D:120:GLU:OE2	1:D:120:GLU:N	2.42	0.49
1:D:141:ILE:O	1:D:141:ILE:HG23	2.12	0.49
1:B:9:TYR:HA	1:B:18:TRP:O	2.13	0.49
1:D:204:LEU:CG	1:D:234:ILE:HD11	2.42	0.49
1:D:40:GLN:HG2	1:D:174:VAL:HG21	1.93	0.49
1:A:260:THR:CG2	1:A:261:SER:N	2.75	0.49
1:C:71:PRO:O	1:C:74:THR:OG1	2.29	0.49
1:D:28:ASP:CA	1:D:55:ARG:O	2.55	0.49
1:C:108:ALA:HB1	1:C:130:VAL:HG22	1.95	0.49
1:D:31:LEU:HB3	1:D:42:PHE:CE1	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:VAL:O	1:A:262:ARG:HB2	2.13	0.48
1:D:172:PRO:HD3	4:D:312:HOH:O	2.12	0.48
1:D:153:MET:HA	1:D:156:ASN:HB2	1.95	0.48
1:C:194:PRO:HB2	1:C:195:PRO:HD3	1.94	0.48
1:D:78:ILE:HG23	1:D:83:LEU:HG	1.94	0.48
1:A:98:ILE:HD12	1:A:121:ARG:C	2.39	0.48
1:D:243:HIS:C	1:D:245:PRO:HD3	2.39	0.48
1:B:44:LYS:HB2	1:B:45:PRO:HD3	1.95	0.47
1:D:124:ASN:ND2	1:D:264:TYR:CD1	2.82	0.47
1:B:5:ARG:NH1	1:B:21:GLU:OE1	2.47	0.47
1:C:263:LYS:HE3	1:C:264:TYR:CZ	2.50	0.47
1:D:248:SER:OG	1:D:249:HIS:N	2.47	0.47
1:C:87:ILE:HD12	1:C:110:THR:HG21	1.96	0.47
1:C:145:ASP:OD1	1:C:148:THR:CB	2.62	0.47
1:D:32:ILE:HG23	1:D:61:MET:HE2	1.96	0.47
1:D:112:LEU:HD21	1:D:127:PRO:HG3	1.95	0.47
1:C:193:ILE:N	1:C:194:PRO:CD	2.78	0.47
1:B:189:TYR:HB2	1:B:190:PRO:HD3	1.97	0.47
1:D:193:ILE:HB	1:D:194:PRO:HD3	1.97	0.47
1:C:142:HIS:NE2	1:C:191:ARG:HD3	2.29	0.47
1:A:44:LYS:N	1:A:45:PRO:HD2	2.30	0.47
1:B:18:TRP:CD2	1:B:90:LEU:HD22	2.50	0.47
1:C:194:PRO:HB2	1:C:195:PRO:CD	2.44	0.46
1:C:232:GLU:OE1	1:C:232:GLU:HA	2.15	0.46
1:D:99:ALA:O	1:D:123:ARG:N	2.39	0.46
1:A:31:LEU:HD23	1:A:102:TRP:HB3	1.96	0.46
1:B:185:TRP:CH2	1:B:190:PRO:HG3	2.51	0.46
1:D:18:TRP:HA	1:D:67:SER:HB3	1.97	0.46
1:D:32:ILE:CG2	1:D:61:MET:HE2	2.46	0.46
1:D:32:ILE:HD12	1:D:101:VAL:HG12	1.97	0.46
1:D:66:ARG:HD3	4:D:302:HOH:O	2.16	0.46
1:D:211:TRP:CE3	1:D:229:ALA:HB2	2.51	0.46
1:A:112:LEU:HD21	1:A:127:PRO:HG3	1.97	0.46
1:C:211:TRP:NE1	1:C:225:ASN:OD1	2.39	0.46
1:D:88:ILE:CD1	1:D:91:LEU:HD12	2.44	0.46
1:C:61:MET:HB3	1:C:62:PRO:CD	2.45	0.46
1:C:108:ALA:CB	1:C:130:VAL:HG22	2.46	0.46
1:C:42:PHE:C	1:C:45:PRO:HD2	2.40	0.45
1:C:54:PHE:CD1	1:C:54:PHE:N	2.84	0.45
1:B:213:VAL:O	1:B:238:THR:HA	2.15	0.45
1:D:33:PRO:HA	1:D:42:PHE:HE2	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:193:ILE:HB	1:C:194:PRO:HD3	1.97	0.45
1:D:221:LEU:HA	1:D:221:LEU:HD23	1.67	0.45
1:B:162:GLY:HA2	1:B:242:ASN:OD1	2.17	0.45
1:D:124:ASN:ND2	1:D:264:TYR:HE1	2.07	0.45
1:A:262:ARG:HD3	1:A:262:ARG:HA	1.76	0.45
1:C:44:LYS:N	1:C:45:PRO:CD	2.79	0.45
1:C:251:GLU:H	1:C:251:GLU:HG2	1.45	0.45
1:B:118:TYR:N	1:B:119:PRO:CD	2.79	0.45
1:B:129:GLU:HB3	1:B:222:PHE:CE2	2.51	0.45
1:D:29:VAL:O	1:D:56:VAL:CG2	2.63	0.45
1:D:239:LEU:HG	1:D:240:PRO:HD2	1.99	0.45
1:C:158:ARG:O	1:C:161:SER:OG	2.32	0.45
1:A:177:ARG:HD3	4:A:453:HOH:O	2.17	0.45
1:D:213:VAL:O	1:D:238:THR:HA	2.17	0.45
1:B:248:SER:HG	1:B:249:HIS:CE1	2.27	0.44
1:C:112:LEU:HA	1:C:112:LEU:HD23	1.75	0.44
1:A:200:LYS:HB2	1:A:203:ASP:OD2	2.17	0.44
1:B:102:TRP:CD1	1:B:102:TRP:C	2.95	0.44
1:D:29:VAL:O	1:D:56:VAL:CA	2.60	0.44
1:B:7:ARG:NH2	4:B:408:HOH:O	2.49	0.44
1:B:105:ALA:HA	1:B:129:GLU:O	2.18	0.44
1:B:63:GLY:C	1:B:68:SER:HA	2.42	0.44
1:D:158:ARG:HD2	1:D:159:ALA:N	2.33	0.44
1:C:141:ILE:HG23	1:C:186:ALA:HB1	1.99	0.44
1:B:89:THR:O	1:B:93:THR:HG23	2.18	0.44
1:B:222:PHE:O	1:B:225:ASN:HB2	2.18	0.43
1:D:204:LEU:CD2	1:D:232:GLU:HG3	2.48	0.43
1:A:45:PRO:O	1:A:49:ILE:HD12	2.18	0.43
1:D:178:LEU:HD23	1:D:178:LEU:HA	1.80	0.43
1:D:225:ASN:N	1:D:225:ASN:HD22	2.16	0.43
1:C:253:PHE:O	1:C:257:VAL:HG23	2.18	0.43
1:C:145:ASP:OD1	1:C:148:THR:OG1	2.23	0.43
1:D:206:LYS:C	1:D:207:VAL:HG13	2.44	0.43
1:A:223:PHE:CE1	1:A:227:VAL:HG11	2.53	0.43
1:D:32:ILE:HG23	1:D:61:MET:CE	2.49	0.43
1:D:131:PRO:O	1:D:194:PRO:HA	2.19	0.43
1:B:104:CYS:HA	1:B:128:HIS:O	2.19	0.43
1:C:46:MET:HG2	1:C:56:VAL:HG11	1.99	0.43
1:D:102:TRP:CD1	1:D:102:TRP:C	2.96	0.43
2:C:301:ZER:O10	2:C:301:ZER:C1P	2.64	0.43
1:A:19:TYR:CE2	1:A:66:ARG:CZ	3.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:5:ARG:HA	1:D:22:GLN:O	2.19	0.42
1:D:160:ALA:HB1	1:D:243:HIS:HB2	2.00	0.42
1:C:78:ILE:HG13	1:C:189:TYR:CE1	2.54	0.42
1:D:42:PHE:O	1:D:46:MET:N	2.52	0.42
1:C:139:LEU:HD22	1:C:140:HIS:CE1	2.53	0.42
1:A:88:ILE:HD13	1:A:88:ILE:HA	1.85	0.42
1:B:166:ALA:HA	1:B:169:ALA:HB3	2.02	0.42
2:A:301:ZER:O10	2:A:301:ZER:H11	2.19	0.42
1:C:4:THR:N	4:C:408:HOH:O	2.51	0.42
1:A:194:PRO:N	1:A:195:PRO:CD	2.82	0.42
1:C:145:ASP:OD1	1:C:148:THR:HB	2.19	0.42
1:A:216:SER:HB2	1:A:242:ASN:OD1	2.20	0.42
1:D:33:PRO:HA	1:D:42:PHE:CE2	2.54	0.42
1:D:246:TYR:HB3	1:D:253:PHE:CE2	2.54	0.42
1:C:39:CYS:HB3	1:C:59:PHE:O	2.20	0.42
1:C:256:TYR:CD1	1:C:256:TYR:C	2.97	0.42
1:C:258:VAL:O	1:C:262:ARG:HB3	2.20	0.41
1:D:102:TRP:HE3	1:D:126:MET:HG3	1.84	0.41
1:A:71:PRO:HA	1:A:72:PRO:HD3	1.84	0.41
1:A:257:VAL:C	1:A:260:THR:HG22	2.45	0.41
1:C:220:LYS:HG2	1:D:215:ALA:C	2.44	0.41
1:A:171:GLY:HA2	1:A:172:PRO:HD3	1.84	0.41
1:D:118:TYR:N	1:D:119:PRO:HD3	2.35	0.41
1:D:39:CYS:HB3	1:D:59:PHE:C	2.45	0.41
1:D:49:ILE:HD11	1:D:258:VAL:HG12	2.02	0.41
1:B:84:ALA:HB1	1:B:114:LEU:HD23	2.03	0.41
1:B:119:PRO:HG3	1:B:206:LYS:HD3	2.02	0.41
1:B:149:ILE:HG21	1:B:183:PRO:HA	2.02	0.41
1:D:153:MET:HE3	1:D:156:ASN:HB2	2.03	0.41
1:A:221:LEU:HD23	1:A:221:LEU:HA	1.92	0.41
1:B:87:ILE:O	1:B:91:LEU:HG	2.21	0.41
1:D:10:VAL:CG1	1:D:93:THR:HG21	2.50	0.41
1:D:72:PRO:HA	1:D:75:TYR:CE1	2.55	0.41
1:C:134:ASN:HD21	1:C:194:PRO:HG3	1.85	0.41
1:B:44:LYS:N	1:B:45:PRO:HD2	2.35	0.41
1:C:71:PRO:HD2	1:C:74:THR:HG21	2.03	0.41
1:B:182:TYR:HB2	1:B:183:PRO:HD3	2.02	0.41
1:D:244:PHE:O	1:D:248:SER:N	2.46	0.41
1:A:129:GLU:HB3	1:A:222:PHE:CZ	2.56	0.40
1:D:49:ILE:HD11	1:D:258:VAL:HG13	1.86	0.40
1:A:126:MET:HE3	1:A:126:MET:HB2	1.80	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:142:HIS:ND1	1:C:143:GLU:HG2	2.37	0.40
1:B:144:VAL:CG2	1:B:149:ILE:HG13	2.48	0.40
1:D:146:PRO:HG3	1:D:187:TYR:OH	2.22	0.40
1:A:90:LEU:HD11	1:A:94:LEU:HD11	2.03	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	261/263 (99%)	253 (97%)	6 (2%)	2 (1%)	16	19
1	B	261/263 (99%)	251 (96%)	8 (3%)	2 (1%)	16	19
1	C	261/263 (99%)	251 (96%)	8 (3%)	2 (1%)	16	19
1	D	259/263 (98%)	232 (90%)	25 (10%)	2 (1%)	16	19
All	All	1042/1052 (99%)	987 (95%)	47 (4%)	8 (1%)	16	19

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	163	ASN
1	D	189	TYR
1	A	189	TYR
1	C	163	ASN
1	C	189	TYR
1	D	164	VAL
1	B	163	ASN
1	B	189	TYR

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	222/222 (100%)	208 (94%)	14 (6%)	16	22
1	B	222/222 (100%)	214 (96%)	8 (4%)	31	45
1	C	222/222 (100%)	201 (90%)	21 (10%)	8	10
1	D	220/222 (99%)	162 (74%)	58 (26%)	0	0
All	All	886/888 (100%)	785 (89%)	101 (11%)	5	6

All (101) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	SER
1	A	51	SER
1	A	89	THR
1	A	93	THR
1	A	95	ASP
1	A	97	LYS
1	A	126	MET
1	A	140	HIS
1	A	151	GLN
1	A	193	ILE
1	A	227	VAL
1	A	255	LYS
1	A	262	ARG
1	A	266	LYS
1	B	10	VAL
1	B	93	THR
1	B	133	GLU
1	B	139	LEU
1	B	164	VAL
1	B	231	ARG
1	B	236	ILE
1	B	255	LYS
1	C	46	MET
1	C	82	LYS

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Mol	Chain	Res	Type
1	C	88	ILE
1	C	95	ASP
1	C	109	SER
1	C	143	GLU
1	C	144	VAL
1	C	173	GLU
1	C	191	ARG
1	C	202	GLU
1	C	204	LEU
1	C	212	THR
1	C	213	VAL
1	C	219	THR
1	C	220	LYS
1	C	225	ASN
1	C	228	ILE
1	C	231	ARG
1	C	246	TYR
1	C	251	GLU
1	C	263	LYS
1	D	5	ARG
1	D	6	THR
1	D	7	ARG
1	D	13	LYS
1	D	29	VAL
1	D	30	VAL
1	D	31	LEU
1	D	44	LYS
1	D	46	MET
1	D	47	SER
1	D	48	LEU
1	D	49	ILE
1	D	55	ARG
1	D	58	THR
1	D	66	ARG
1	D	67	SER
1	D	68	SER
1	D	74	THR
1	D	78	ILE
1	D	87	ILE
1	D	88	ILE
1	D	90	LEU
1	D	97	LYS

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Mol	Chain	Res	Type
1	D	98	ILE
1	D	101	VAL
1	D	106	SER
1	D	109	SER
1	D	118	TYR
1	D	120	GLU
1	D	121	ARG
1	D	122	VAL
1	D	123	ARG
1	D	133	GLU
1	D	138	LEU
1	D	139	LEU
1	D	141	ILE
1	D	143	GLU
1	D	157	SER
1	D	158	ARG
1	D	164	VAL
1	D	173	GLU
1	D	184	ARG
1	D	192	THR
1	D	202	GLU
1	D	204	LEU
1	D	206	LYS
1	D	209	ILE
1	D	211	TRP
1	D	212	THR
1	D	217	THR
1	D	224	GLU
1	D	225	ASN
1	D	226	ILE
1	D	227	VAL
1	D	239	LEU
1	D	262	ARG
1	D	263	LYS
1	D	264	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	134	ASN
1	A	151	GLN

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Mol	Chain	Res	Type
1	C	134	ASN
1	C	205	HIS
1	D	22	GLN
1	D	124	ASN
1	D	156	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	ZER	B	301	-	24,24,24	1.64	5 (20%)	32,32,32	1.11	3 (9%)
2	ZER	A	301	-	24,24,24	1.62	5 (20%)	32,32,32	1.12	2 (6%)
3	SO4	A	302	-	4,4,4	0.50	0	6,6,6	0.24	0
2	ZER	C	301	-	24,24,24	1.61	5 (20%)	32,32,32	0.91	0

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	ZER	B	301	-	-	7/22/22/22	0/1/2/2
2	ZER	A	301	-	-	7/22/22/22	0/1/2/2
2	ZER	C	301	-	-	8/22/22/22	0/1/2/2

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	ZER	C1-C12	-3.69	1.42	1.50
2	A	301	ZER	C6-C1P	-3.68	1.41	1.47
2	A	301	ZER	C1-C12	-3.48	1.42	1.50
2	B	301	ZER	C5P-C6P	-3.47	1.41	1.51
2	C	301	ZER	C5P-C6P	-3.47	1.41	1.51
2	B	301	ZER	C1-C12	-3.41	1.42	1.50
2	B	301	ZER	C1P-C2P	3.33	1.41	1.31
2	A	301	ZER	C5P-C6P	-3.32	1.41	1.51
2	A	301	ZER	C1P-C2P	3.16	1.41	1.31
2	B	301	ZER	C7P-C6P	-3.10	1.42	1.51
2	B	301	ZER	C6-C1P	-3.09	1.42	1.47
2	C	301	ZER	C6-C1P	-3.05	1.42	1.47
2	C	301	ZER	C7P-C6P	-2.89	1.42	1.51
2	C	301	ZER	C1P-C2P	2.71	1.40	1.31
2	A	301	ZER	C7P-C6P	-2.60	1.43	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	ZER	C3P-C2P-C1P	-2.64	119.63	125.48
2	B	301	ZER	C7P-C6P-C5P	2.61	121.89	116.95
2	A	301	ZER	C11-C10-C9P	-2.56	107.28	113.90
2	B	301	ZER	C8P-C7P-C6P	-2.29	108.78	114.59
2	A	301	ZER	O2-C2-C1	-2.03	117.38	121.14

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	ZER	C1-C12-O10-C10
2	C	301	ZER	C3P-C4P-C5P-C6P
2	B	301	ZER	C2P-C3P-C4P-C5P
2	A	301	ZER	C3P-C4P-C5P-C6P

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Mol	Chain	Res	Type	Atoms
2	B	301	ZER	C3P-C4P-C5P-C6P
2	B	301	ZER	C6P-C7P-C8P-C9P
2	C	301	ZER	C6P-C7P-C8P-C9P
2	C	301	ZER	C2P-C3P-C4P-C5P
2	A	301	ZER	C2P-C3P-C4P-C5P
2	A	301	ZER	O12-C12-O10-C10
2	A	301	ZER	C6P-C7P-C8P-C9P
2	C	301	ZER	C5P-C6P-C7P-C8P
2	C	301	ZER	C11-C10-O10-C12
2	C	301	ZER	C1P-C2P-C3P-C4P
2	C	301	ZER	O6P-C6P-C7P-C8P
2	B	301	ZER	C2-C1-C12-O10
2	C	301	ZER	C9P-C10-O10-C12
2	B	301	ZER	C1-C12-O10-C10
2	A	301	ZER	O6P-C6P-C7P-C8P
2	B	301	ZER	O6P-C6P-C7P-C8P
2	B	301	ZER	C6-C1-C12-O10
2	A	301	ZER	C7P-C8P-C9P-C10

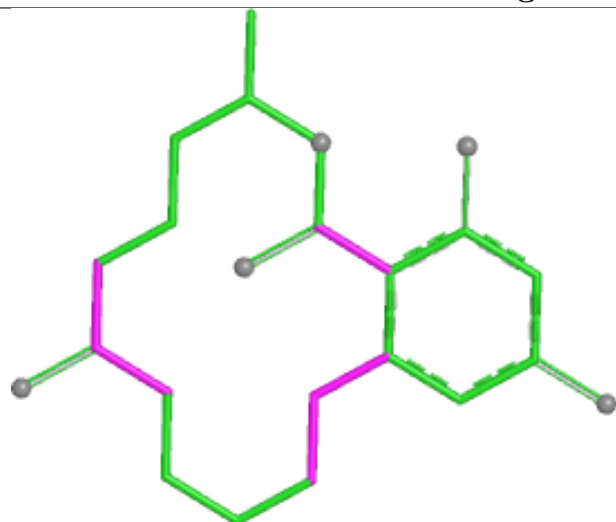
There are no ring outliers.

2 monomers are involved in 6 short contacts:

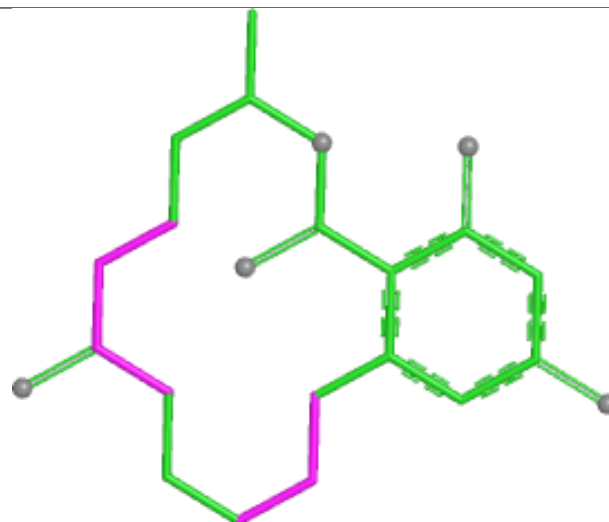
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	ZER	3	0
2	C	301	ZER	3	0

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.

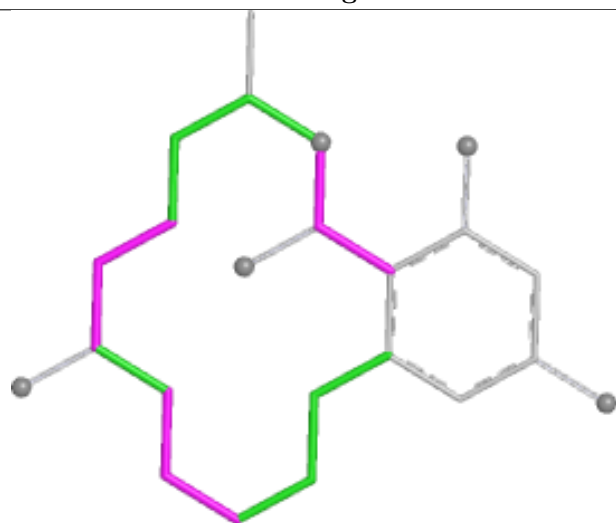
Ligand ZER B 301



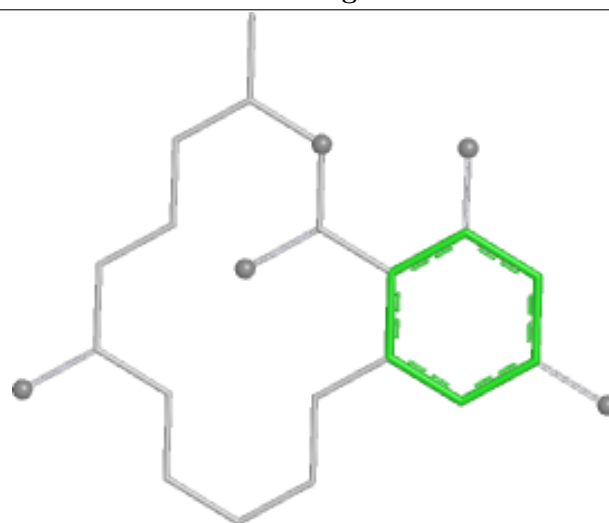
Bond lengths



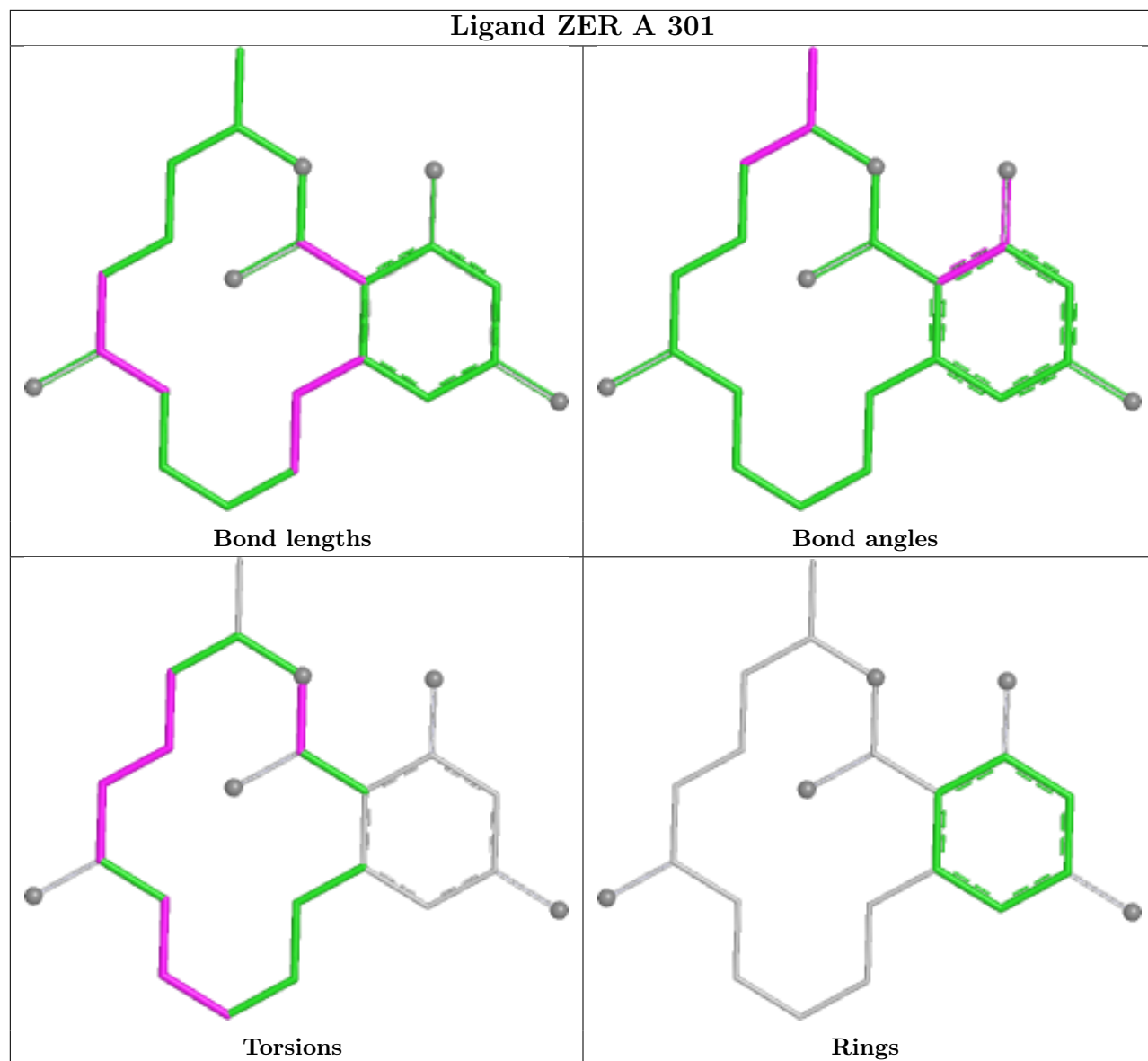
Bond angles

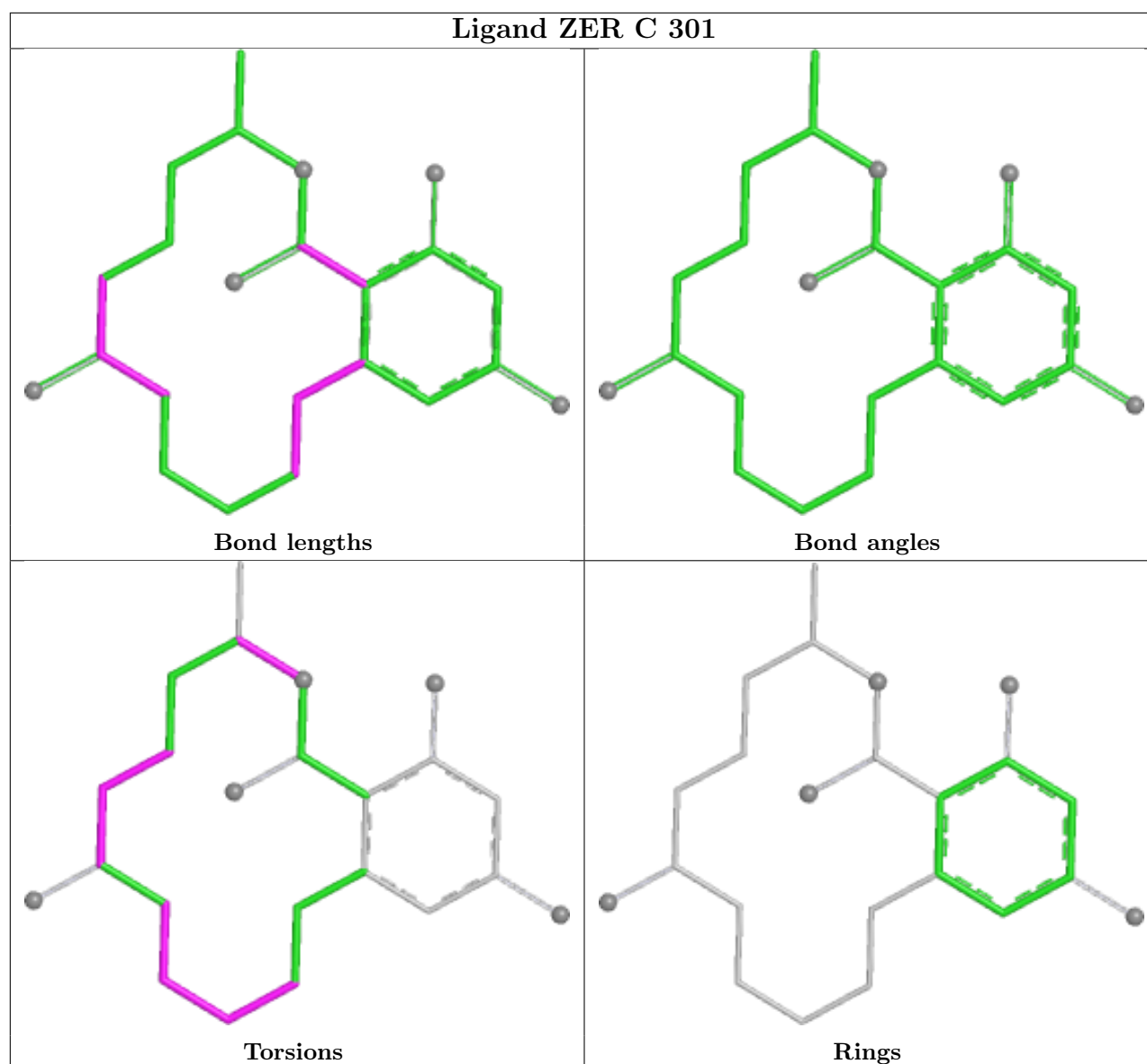


Torsions



Rings





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	263/263 (100%)	-1.50	0 100 100	23, 36, 50, 88	0
1	B	263/263 (100%)	-1.52	0 100 100	22, 35, 49, 76	0
1	C	263/263 (100%)	-1.41	0 100 100	31, 48, 65, 81	0
1	D	261/263 (99%)	-1.08	0 100 100	34, 70, 100, 113	0
All	All	1050/1052 (99%)	-1.38	0 100 100	22, 44, 85, 113	0

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](#)

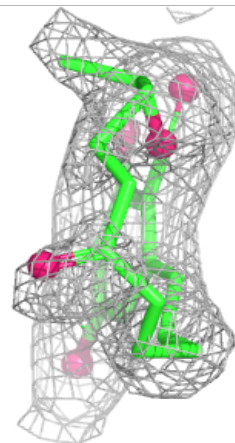
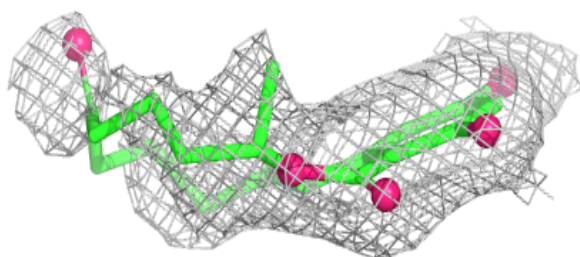
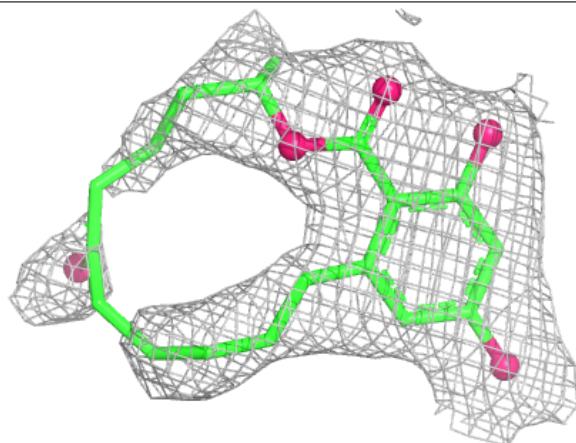
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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	ZER	A	301	23/23	0.99	0.03	22,33,58,62	0
2	ZER	B	301	23/23	0.99	0.03	20,32,45,60	0
2	ZER	C	301	23/23	0.99	0.04	33,45,72,86	0
3	SO4	A	302	5/5	0.99	0.05	48,55,75,76	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.

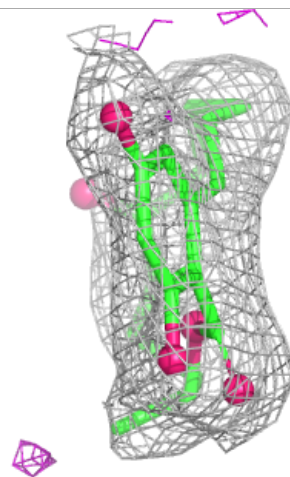
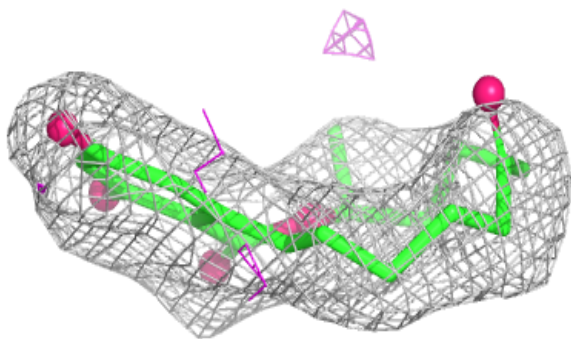
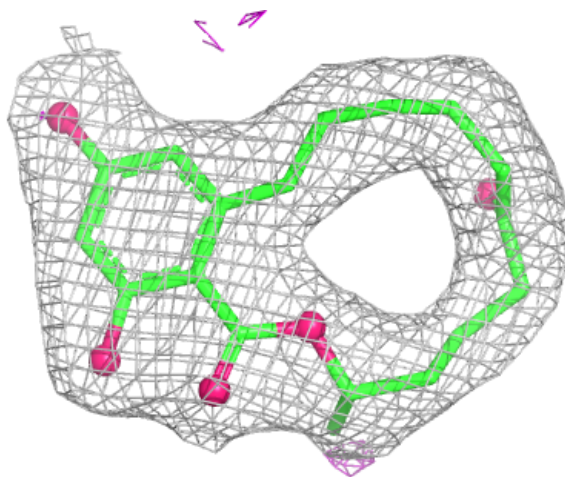
Electron density around ZER A 301:

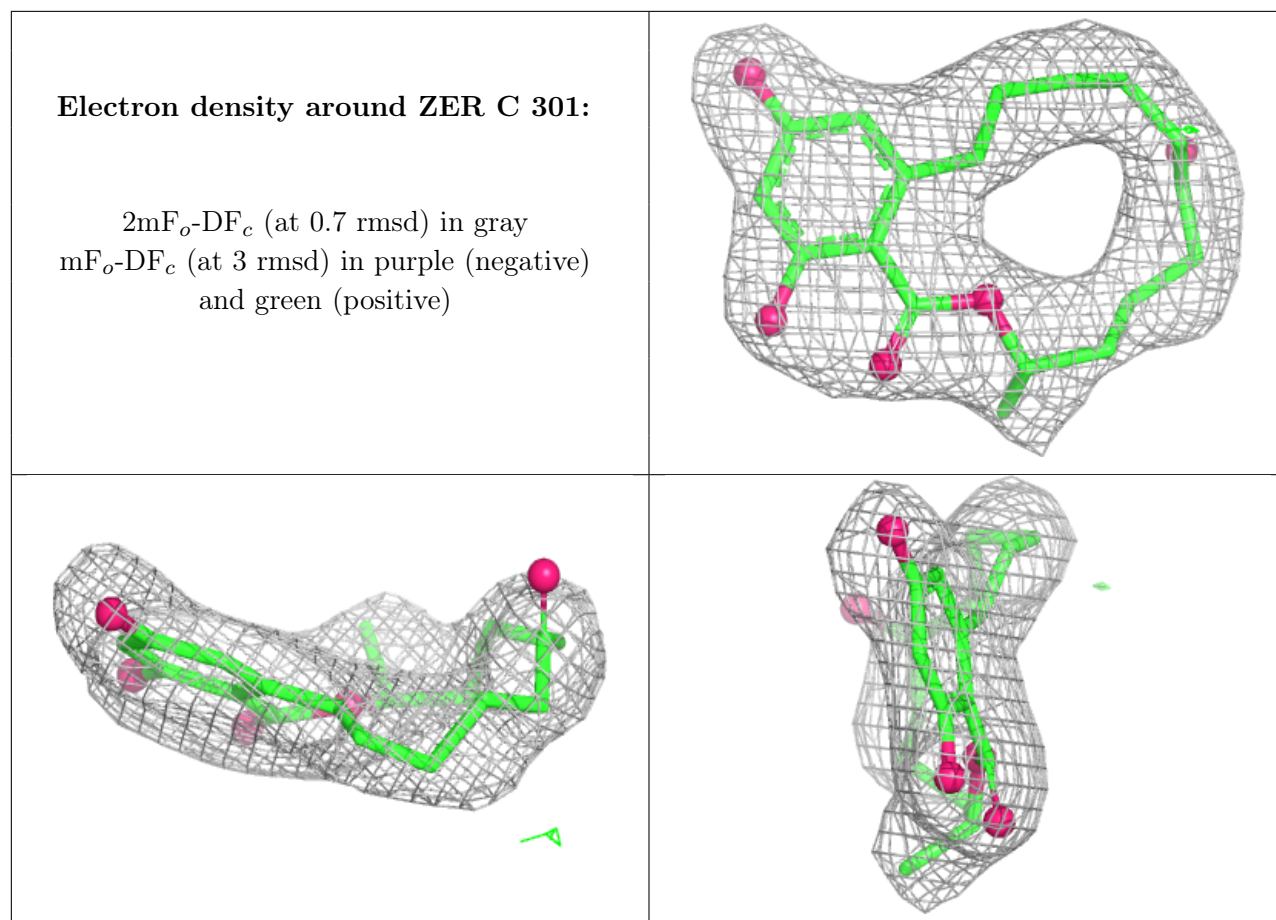
2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ZER B 301:

$2mF_o-DF_c$ (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.