



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:30 PM UTC

PDB ID : 2ZX7 / pdb_00002zx7
Title : alpha-L-fucosidase complexed with inhibitor, F10-2C
Authors : Wu, H.-J.; Ko, T.-P.; Ho, C.-W.; Lin, C.-H.; Wang, A.H.-J.
Deposited on : 2008-12-19
Resolution : 2.48 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

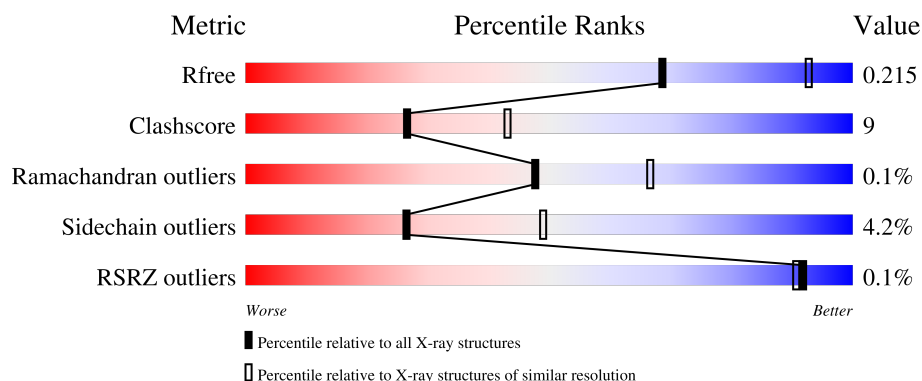
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.48 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	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (2.50-2.46)

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	455	 71% 22% . .
1	B	455	 72% 22% . .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7786 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.

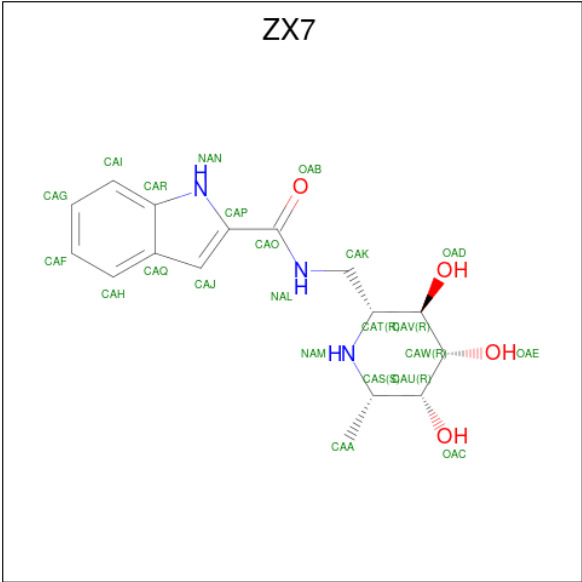
- Molecule 1 is a protein called Alpha-L-fucosidase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	442	Total	C	N	O	S	0	0	0
			3648	2377	600	663	8			
1	B	442	Total	C	N	O	S	0	0	0
			3648	2377	600	663	8			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	450	HIS	-	expression tag	UNP Q9WYE2
A	451	HIS	-	expression tag	UNP Q9WYE2
A	452	HIS	-	expression tag	UNP Q9WYE2
A	453	HIS	-	expression tag	UNP Q9WYE2
A	454	HIS	-	expression tag	UNP Q9WYE2
A	455	HIS	-	expression tag	UNP Q9WYE2
B	450	HIS	-	expression tag	UNP Q9WYE2
B	451	HIS	-	expression tag	UNP Q9WYE2
B	452	HIS	-	expression tag	UNP Q9WYE2
B	453	HIS	-	expression tag	UNP Q9WYE2
B	454	HIS	-	expression tag	UNP Q9WYE2
B	455	HIS	-	expression tag	UNP Q9WYE2

- Molecule 2 is N-[[[(2R,3R,4R,5R,6S)-3,4,5-trihydroxy-6-methylpiperidin-2-yl]methyl}-1H-indole-2-carboxamide (CCD ID: ZX7) (formula: C₁₆H₂₁N₃O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			23	16	3	4		
2	B	1	Total	C	N	O	0	0
			23	16	3	4		

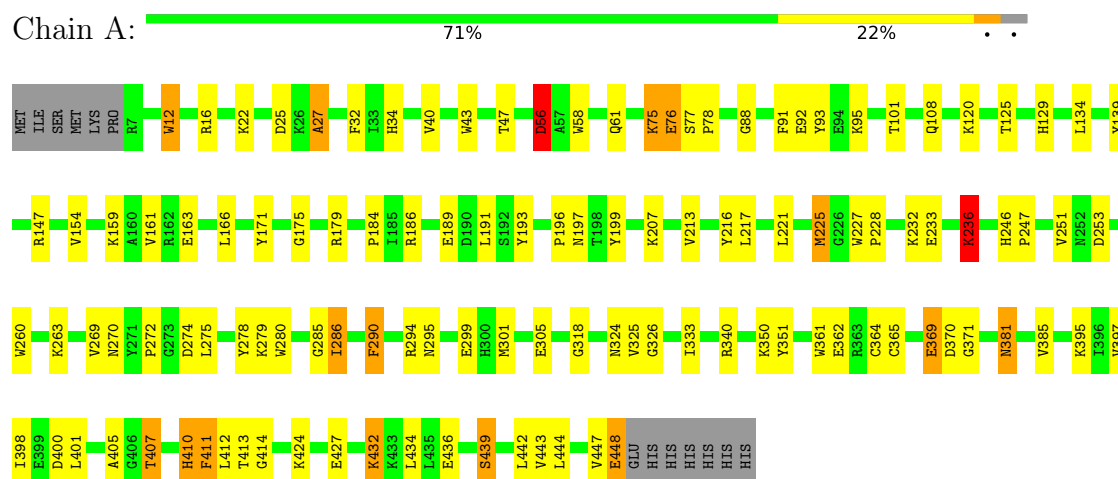
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	214	Total	O	0	0
			214	214		
3	B	230	Total	O	0	0
			230	230		

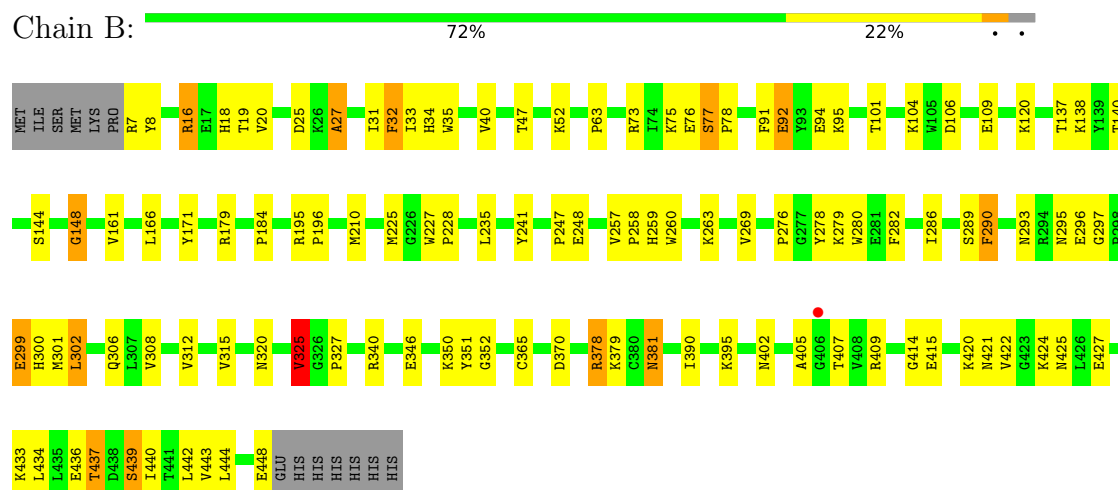
3 Residue-property plots

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: Alpha-L-fucosidase, putative



- Molecule 1: Alpha-L-fucosidase, putative



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	180.39Å 180.39Å 169.31Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.48 30.00 – 2.48	Depositor EDS
% Data completeness (in resolution range)	95.9 (30.00-2.48) 95.8 (30.00-2.48)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.02 (at 2.48Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.161 , 0.216 0.162 , 0.215	Depositor DCC
R_{free} test set	1784 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	35.8	Xtriage
Anisotropy	0.151	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 60.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7786	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.15% 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: ZX7

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	1.12	3/3768 (0.1%)	1.26	33/5120 (0.6%)
1	B	1.13	6/3768 (0.2%)	1.25	26/5120 (0.5%)
All	All	1.12	9/7536 (0.1%)	1.25	59/10240 (0.6%)

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	A	0	2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	213	VAL	CA-CB	8.76	1.64	1.54
1	B	31	ILE	CA-CB	8.05	1.63	1.54
1	A	225	MET	SD-CE	-6.74	1.62	1.79
1	B	225	MET	SD-CE	-6.44	1.63	1.79
1	B	315	VAL	CA-CB	6.16	1.62	1.54
1	A	101	THR	CA-CB	5.97	1.62	1.54
1	B	301	MET	SD-CE	5.55	1.93	1.79
1	B	33	ILE	C-O	-5.34	1.18	1.24
1	B	325	VAL	CA-CB	5.20	1.61	1.54

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	77	SER	N-CA-C	-10.23	92.90	109.58
1	B	77	SER	N-CA-C	-9.20	94.37	109.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	286	ILE	N-CA-C	-8.83	101.80	110.72
1	B	365	CYS	N-CA-C	8.18	121.91	109.23
1	A	290	PHE	N-CA-C	-7.42	102.82	112.68
1	B	440	ILE	N-CA-C	-7.38	106.69	113.71
1	B	40	VAL	CA-C-N	-7.28	111.27	119.28
1	B	40	VAL	C-N-CA	-7.28	111.27	119.28
1	A	199	TYR	N-CA-C	-7.27	103.29	111.14
1	A	175	GLY	N-CA-C	7.23	123.06	113.37
1	A	365	CYS	N-CA-C	7.07	120.19	109.23
1	A	371	GLY	N-CA-C	7.07	125.40	115.43
1	A	411	PHE	N-CA-C	6.96	118.87	111.28
1	B	352	GLY	N-CA-C	6.76	122.26	113.27
1	B	148	GLY	N-CA-C	6.74	126.09	112.34
1	A	75	LYS	CA-C-N	-6.70	113.03	122.34
1	A	75	LYS	C-N-CA	-6.70	113.03	122.34
1	B	415	GLU	N-CA-C	6.61	119.81	109.96
1	B	144	SER	N-CA-C	6.58	120.76	112.87
1	A	326	GLY	CA-C-N	6.56	126.51	120.21
1	A	326	GLY	C-N-CA	6.56	126.51	120.21
1	B	290	PHE	N-CA-C	-6.56	103.96	112.68
1	A	56	ASP	N-CA-C	-6.51	104.81	112.89
1	A	40	VAL	CA-C-N	-6.48	112.33	119.19
1	A	40	VAL	C-N-CA	-6.48	112.33	119.19
1	B	101	THR	N-CA-C	6.37	121.20	113.17
1	A	318	GLY	CA-C-N	-6.36	115.75	122.30
1	A	318	GLY	C-N-CA	-6.36	115.75	122.30
1	B	269	VAL	N-CA-C	6.29	117.14	108.84
1	A	47	THR	N-CA-C	6.24	118.08	111.28
1	B	47	THR	N-CA-C	6.13	117.96	111.28
1	B	320	ASN	N-CA-C	-6.11	99.95	109.72
1	A	129	HIS	N-CA-C	5.96	119.37	111.75
1	A	286	ILE	N-CA-C	-5.92	103.70	111.09
1	A	236	LYS	N-CA-C	-5.80	104.96	111.28
1	A	407	THR	N-CA-C	5.79	117.87	108.26
1	B	442	LEU	N-CA-C	-5.65	100.78	109.76
1	B	302	LEU	N-CA-C	-5.64	102.13	110.48
1	A	442	LEU	N-CA-C	-5.62	100.65	109.25
1	A	439	SER	N-CA-C	5.61	120.25	112.45
1	A	269	VAL	N-CA-C	5.56	116.18	108.84
1	B	282	PHE	CA-C-N	-5.56	113.94	122.62
1	B	282	PHE	C-N-CA	-5.56	113.94	122.62
1	A	410	HIS	N-CA-C	-5.51	100.20	108.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	439	SER	CA-C-N	-5.51	113.82	121.53
1	B	439	SER	C-N-CA	-5.51	113.82	121.53
1	A	88	GLY	N-CA-C	5.50	119.55	112.18
1	A	232	LYS	N-CA-C	5.48	117.25	111.28
1	B	179	ARG	N-CA-C	-5.47	106.56	113.18
1	B	27	ALA	N-CA-C	5.41	118.09	111.82
1	A	27	ALA	N-CA-C	5.39	117.23	111.36
1	A	414	GLY	N-CA-C	-5.38	107.92	114.92
1	B	351	TYR	N-CA-C	5.36	119.92	113.17
1	B	259	HIS	N-CA-C	5.35	118.83	110.32
1	A	58	TRP	N-CA-C	5.28	120.08	111.37
1	A	12	TRP	N-CA-C	5.22	116.97	111.28
1	A	179	ARG	N-CA-C	-5.15	106.95	113.18
1	B	32	PHE	N-CA-C	-5.13	102.17	110.32
1	A	225	MET	N-CA-C	5.03	118.72	112.58

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	193	TYR	Sidechain
1	A	93	TYR	Sidechain

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	3648	0	3512	78	0
1	B	3648	0	3512	63	0
2	A	23	0	21	2	0
2	B	23	0	21	1	0
3	A	214	0	0	1	1
3	B	230	0	0	3	0
All	All	7786	0	7066	134	1

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 (134) 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:A:395:LYS:HE3	1:A:427:GLU:OE1	1.66	0.95
1:A:16:ARG:HD2	1:B:16:ARG:HH21	1.31	0.94
1:A:381:ASN:HD22	1:A:381:ASN:H	1.25	0.83
1:B:297:GLY:H	1:B:300:HIS:HD2	1.26	0.82
1:A:260:TRP:CZ2	1:B:16:ARG:HG2	2.17	0.79
1:A:381:ASN:HD22	1:A:381:ASN:N	1.83	0.75
1:A:260:TRP:CE2	1:B:16:ARG:HG2	2.22	0.75
1:A:270:ASN:ND2	2:A:901:ZX7:HAG	2.01	0.75
1:A:16:ARG:HH11	1:B:16:ARG:HE	1.34	0.74
1:B:297:GLY:H	1:B:300:HIS:CD2	2.10	0.70
1:A:108:GLN:HE22	1:A:163:GLU:HB2	1.56	0.70
1:B:94:GLU:H	1:B:94:GLU:CD	2.00	0.68
1:A:405:ALA:HA	1:A:448:GLU:HA	1.76	0.67
1:A:290:PHE:HA	1:A:324:ASN:ND2	2.10	0.67
1:A:381:ASN:H	1:A:381:ASN:ND2	1.93	0.66
1:A:285:GLY:HA2	1:A:324:ASN:HB3	1.78	0.66
1:A:286:ILE:CG2	1:A:301:MET:HE3	2.25	0.66
1:B:161:VAL:HG13	1:B:166:LEU:HB2	1.79	0.64
1:A:286:ILE:HG22	1:A:301:MET:HE3	1.79	0.64
1:A:196:PRO:O	1:A:197:ASN:HB2	1.99	0.62
1:B:92:GLU:CB	1:B:94:GLU:OE2	2.47	0.62
1:B:263:LYS:HD3	1:B:278:TYR:CE1	2.36	0.61
1:B:91:PHE:CE1	1:B:95:LYS:HD2	2.35	0.61
1:B:409:ARG:NH1	1:B:414:GLY:O	2.34	0.60
1:A:413:THR:HG21	1:A:434:LEU:HD22	1.82	0.60
1:B:92:GLU:HB3	1:B:94:GLU:OE2	2.01	0.60
1:B:106:ASP:OD2	1:B:109:GLU:HG3	2.03	0.59
1:B:227:TRP:CG	1:B:228:PRO:HD2	2.38	0.59
2:B:902:ZX7:HAJ	3:B:600:HOH:O	2.04	0.57
1:A:361:TRP:CG	1:A:362:GLU:H	2.23	0.57
1:B:171:TYR:CD2	1:B:171:TYR:C	2.83	0.56
1:B:434:LEU:O	1:B:437:THR:HG22	2.07	0.56
1:A:405:ALA:HB1	1:A:447:VAL:O	2.06	0.55
1:A:263:LYS:HD3	1:A:278:TYR:CE1	2.41	0.55
1:A:171:TYR:CD2	1:A:171:TYR:C	2.85	0.55
1:A:43:TRP:NE1	1:A:295:ASN:HD21	2.05	0.55
1:A:301:MET:HE2	1:A:340:ARG:HD2	1.90	0.54
1:B:16:ARG:HG3	1:B:16:ARG:O	2.08	0.54
1:B:34:HIS:CG	1:B:290:PHE:HB3	2.44	0.53
1:A:75:LYS:O	1:A:76:GLU:CB	2.55	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:TYR:CE2	1:A:207:LYS:HE2	2.44	0.53
1:B:395:LYS:HE2	1:B:427:GLU:CD	2.35	0.52
1:B:378:ARG:CG	1:B:378:ARG:HH11	2.21	0.52
1:A:233:GLU:OE1	1:A:236:LYS:HE3	2.10	0.52
1:A:34:HIS:CG	1:A:290:PHE:HB3	2.44	0.52
1:A:325:VAL:HG11	1:A:333:ILE:HD13	1.92	0.52
1:B:433:LYS:HG3	1:B:434:LEU:N	2.25	0.51
1:A:184:PRO:HG2	1:A:186:ARG:NH2	2.25	0.51
1:B:35:TRP:CD1	1:B:327:PRO:HB2	2.46	0.51
1:B:350:LYS:HE2	1:B:439:SER:HB2	1.93	0.51
1:B:378:ARG:HG2	1:B:378:ARG:NH1	2.25	0.51
1:A:108:GLN:NE2	1:A:163:GLU:HB2	2.24	0.50
1:A:161:VAL:HG13	1:A:166:LEU:HB2	1.94	0.50
1:A:221:LEU:HB2	1:A:251:VAL:HG12	1.94	0.50
1:A:350:LYS:NZ	1:A:439:SER:HB2	2.26	0.50
1:B:210:MET:HE1	1:B:241:TYR:CE2	2.46	0.50
1:B:297:GLY:N	1:B:300:HIS:HD2	2.02	0.49
1:B:299:GLU:H	1:B:299:GLU:CD	2.21	0.49
1:A:270:ASN:HD21	2:A:901:ZX7:HAG	1.74	0.49
1:A:301:MET:CE	1:A:340:ARG:HD2	2.42	0.49
1:A:125:THR:HA	1:A:171:TYR:HB3	1.95	0.48
1:B:92:GLU:HB2	1:B:94:GLU:OE2	2.13	0.48
1:B:381:ASN:O	1:B:448:GLU:HG2	2.13	0.48
1:A:16:ARG:CG	1:B:260:TRP:CE2	2.96	0.48
1:A:12:TRP:CZ2	1:A:236:LYS:HG2	2.48	0.48
1:A:16:ARG:HG3	1:B:260:TRP:CZ2	2.49	0.48
1:B:296:GLU:HA	1:B:300:HIS:CD2	2.49	0.47
1:B:91:PHE:CD1	1:B:95:LYS:HD2	2.50	0.47
1:A:56:ASP:OD1	3:A:634:HOH:O	2.20	0.47
1:B:137:THR:HG21	1:B:140:THR:HG22	1.96	0.47
1:A:25:ASP:HA	1:A:279:LYS:HE2	1.97	0.47
1:A:91:PHE:CE2	1:A:95:LYS:HB3	2.50	0.47
1:A:16:ARG:HG2	1:B:260:TRP:CE2	2.49	0.46
1:B:27:ALA:HB1	1:B:120:LYS:HB2	1.96	0.46
1:B:161:VAL:CG1	1:B:166:LEU:HB2	2.43	0.46
1:A:227:TRP:CG	1:A:228:PRO:HD2	2.50	0.46
1:A:351:TYR:CD1	1:A:443:VAL:HG21	2.51	0.46
1:A:301:MET:HE2	1:A:340:ARG:CD	2.46	0.46
1:B:210:MET:HE1	1:B:241:TYR:CD2	2.51	0.46
1:B:420:LYS:HE3	1:B:422:VAL:HG22	1.97	0.46
1:A:400:ASP:HA	1:A:424:LYS:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:308:VAL:O	1:B:312:VAL:HG23	2.15	0.46
1:A:159:LYS:O	1:A:163:GLU:HG2	2.16	0.46
1:A:410:HIS:CE1	1:A:412:LEU:HB3	2.50	0.46
1:B:7:ARG:HD3	3:B:622:HOH:O	2.16	0.46
1:A:75:LYS:O	1:A:76:GLU:HB2	2.15	0.45
1:B:302:LEU:HD22	1:B:306:GLN:HB3	1.98	0.45
1:A:385:VAL:HB	1:A:444:LEU:HB2	1.98	0.45
1:A:398:ILE:HG21	1:A:401:LEU:HD23	1.99	0.45
1:B:63:PRO:O	1:B:289:SER:HB2	2.16	0.45
1:A:134:LEU:O	1:A:154:VAL:HG23	2.16	0.45
1:A:432:LYS:O	1:A:436:GLU:HG3	2.16	0.45
1:A:253:ASP:CG	1:A:263:LYS:HA	2.41	0.45
1:A:290:PHE:O	1:A:324:ASN:ND2	2.46	0.45
1:B:293:ASN:OD1	1:B:295:ASN:HB2	2.17	0.45
1:B:402:ASN:HA	1:B:421:ASN:OD1	2.17	0.44
1:B:19:THR:O	1:B:20:VAL:C	2.60	0.44
1:A:305:GLU:CD	1:A:305:GLU:H	2.26	0.44
1:A:411:PHE:CD1	1:A:411:PHE:C	2.95	0.44
1:B:378:ARG:HG3	1:B:379:LYS:N	2.33	0.44
1:B:75:LYS:O	1:B:76:GLU:HB2	2.18	0.44
1:A:61:GLN:HE21	1:A:61:GLN:HB3	1.70	0.44
1:A:217:LEU:HD11	1:A:246:HIS:HB2	1.99	0.44
1:B:8:TYR:CE1	1:B:18:HIS:HB2	2.53	0.44
1:B:325:VAL:O	1:B:327:PRO:HD3	2.18	0.44
1:A:91:PHE:C	1:A:91:PHE:CD2	2.96	0.43
1:A:405:ALA:CA	1:A:448:GLU:HA	2.47	0.43
1:B:340:ARG:HH11	1:B:340:ARG:HG3	1.83	0.43
1:A:22:LYS:HB2	1:A:22:LYS:HE3	1.72	0.43
1:A:444:LEU:HD23	1:A:444:LEU:HA	1.77	0.43
1:B:77:SER:O	1:B:78:PRO:C	2.59	0.43
1:A:191:LEU:HD22	1:A:225:MET:HE1	1.99	0.43
1:A:27:ALA:HB1	1:A:120:LYS:HB2	1.99	0.42
1:B:235:LEU:HD12	1:B:257:VAL:HG11	2.01	0.42
1:B:405:ALA:HA	1:B:448:GLU:HA	2.00	0.42
1:A:216:TYR:O	1:A:217:LEU:C	2.62	0.42
1:A:370:ASP:OD2	1:A:370:ASP:C	2.63	0.42
1:A:246:HIS:O	1:A:247:PRO:C	2.62	0.42
1:A:286:ILE:HG21	1:A:301:MET:HE3	1.98	0.42
1:A:369:GLU:HB2	1:A:397:VAL:HB	2.01	0.41
1:B:73:ARG:HB2	1:B:184:PRO:HB3	2.00	0.41
1:A:274:ASP:CG	1:A:275:LEU:H	2.28	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:424:LYS:HG3	1:B:425:ASN:OD1	2.20	0.41
1:A:275:LEU:CD2	1:A:364:CYS:HB3	2.51	0.41
1:A:189:GLU:H	1:A:189:GLU:CD	2.29	0.41
1:A:272:PRO:CG	1:A:280:TRP:CZ2	3.04	0.41
1:B:443:VAL:O	1:B:444:LEU:HD23	2.21	0.41
1:A:16:ARG:HG2	1:B:260:TRP:NE1	2.36	0.41
1:B:25:ASP:HA	1:B:279:LYS:HE2	2.03	0.41
1:A:290:PHE:CA	1:A:324:ASN:ND2	2.82	0.41
1:B:247:PRO:HD2	3:B:467:HOH:O	2.20	0.40
1:A:272:PRO:HG3	1:A:280:TRP:CZ2	2.56	0.40
1:B:195:ARG:O	1:B:196:PRO:C	2.64	0.40
1:B:276:PRO:HG2	1:B:280:TRP:CD1	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:520:HOH:O	3:A:520:HOH:O[6_555]	1.89	0.31

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	440/455 (97%)	418 (95%)	22 (5%)	0	100	100
1	B	440/455 (97%)	417 (95%)	22 (5%)	1 (0%)	43	61
All	All	880/910 (97%)	835 (95%)	44 (5%)	1 (0%)	48	66

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	148	GLY

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	382/395 (97%)	368 (96%)	14 (4%)	30	54
1	B	382/395 (97%)	364 (95%)	18 (5%)	23	44
All	All	764/790 (97%)	732 (96%)	32 (4%)	26	49

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	PHE
1	A	56	ASP
1	A	76	GLU
1	A	78	PRO
1	A	92	GLU
1	A	147	ARG
1	A	236	LYS
1	A	294	ARG
1	A	299	GLU
1	A	369	GLU
1	A	381	ASN
1	A	407	THR
1	A	432	LYS
1	A	448	GLU
1	B	16	ARG
1	B	32	PHE
1	B	52	LYS
1	B	92	GLU
1	B	104	LYS
1	B	138	LYS
1	B	248	GLU
1	B	258	PRO
1	B	299	GLU
1	B	325	VAL
1	B	346	GLU
1	B	370	ASP
1	B	378	ARG

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Mol	Chain	Res	Type
1	B	381	ASN
1	B	390	ILE
1	B	407	THR
1	B	436	GLU
1	B	437	THR

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	61	GLN
1	A	108	GLN
1	A	208	GLN
1	A	270	ASN
1	A	295	ASN
1	A	381	ASN
1	B	300	HIS

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](#)

2 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	ZX7	B	902	-	25,25,25	1.43	6 (24%)	34,36,36	1.15	4 (11%)
2	ZX7	A	901	-	25,25,25	1.46	5 (20%)	34,36,36	1.33	7 (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	ZX7	B	902	-	-	2/9/29/29	0/3/3/3
2	ZX7	A	901	-	-	4/9/29/29	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	ZX7	CAP-NAN	-3.53	1.33	1.38
2	B	902	ZX7	CAP-NAN	-3.48	1.33	1.38
2	A	901	ZX7	CAQ-CAJ	-3.45	1.33	1.44
2	B	902	ZX7	CAQ-CAJ	-3.36	1.33	1.44
2	A	901	ZX7	CAR-NAN	-2.93	1.33	1.38
2	A	901	ZX7	CAJ-CAP	-2.88	1.33	1.37
2	B	902	ZX7	CAJ-CAP	-2.75	1.33	1.37
2	B	902	ZX7	CAR-NAN	-2.74	1.33	1.38
2	B	902	ZX7	CAP-CAO	2.35	1.53	1.48
2	A	901	ZX7	CAP-CAO	2.26	1.53	1.48
2	B	902	ZX7	CAS-NAM	2.00	1.47	1.45

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	ZX7	CAI-CAR-CAQ	-3.01	119.40	122.13
2	B	902	ZX7	CAI-CAR-CAQ	-2.97	119.44	122.13
2	B	902	ZX7	CAJ-CAP-CAO	-2.91	120.74	130.55
2	A	901	ZX7	CAJ-CAP-CAO	-2.72	121.37	130.55
2	A	901	ZX7	CAR-NAN-CAP	-2.59	106.23	109.01
2	A	901	ZX7	CAJ-CAP-NAN	2.44	111.57	109.15
2	B	902	ZX7	CAR-NAN-CAP	-2.32	106.53	109.01
2	B	902	ZX7	CAJ-CAP-NAN	2.27	111.41	109.15
2	A	901	ZX7	CAO-CAP-NAN	2.20	127.06	120.04
2	A	901	ZX7	CAK-NAL-CAO	-2.12	118.86	122.06
2	A	901	ZX7	OAB-CAO-CAP	-2.07	118.43	121.09

There are no chirality outliers.

All (6) torsion outliers are listed below:

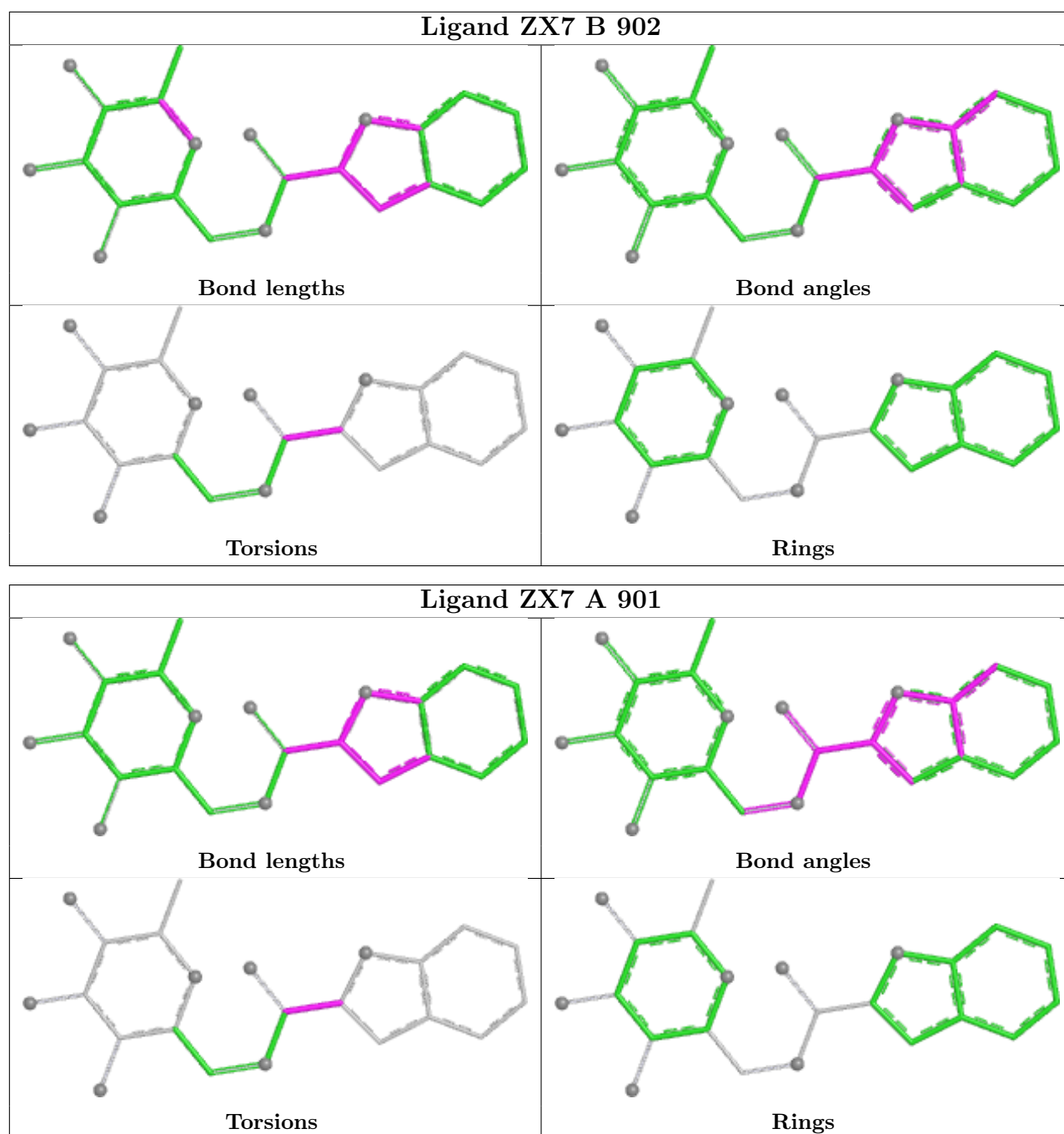
Mol	Chain	Res	Type	Atoms
2	A	901	ZX7	OAB-CAO-CAP-CAJ
2	A	901	ZX7	OAB-CAO-CAP-NAN
2	A	901	ZX7	NAL-CAO-CAP-CAJ
2	A	901	ZX7	NAL-CAO-CAP-NAN
2	B	902	ZX7	NAL-CAO-CAP-CAJ
2	B	902	ZX7	OAB-CAO-CAP-CAJ

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	902	ZX7	1	0
2	A	901	ZX7	2	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	442/455 (97%)	-0.71	0	100 100	18, 34, 59, 80	0
1	B	442/455 (97%)	-0.66	1 (0%)	91 90	21, 35, 61, 81	0
All	All	884/910 (97%)	-0.68	1 (0%)	92 91	18, 35, 60, 81	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	406	GLY	2.4

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.

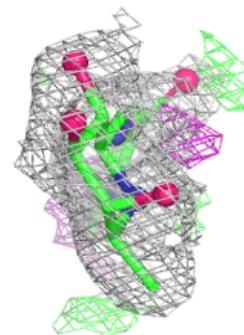
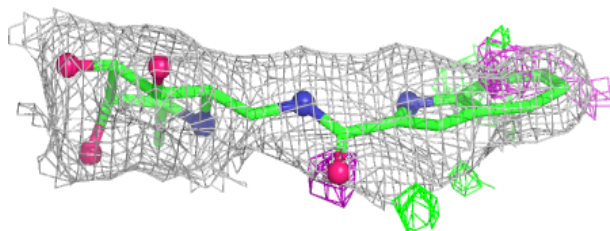
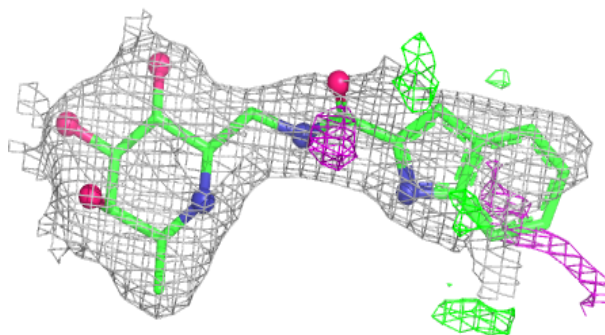
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q < 0.9
2	ZX7	A	901	23/23	0.91	0.11	29,41,63,64	0
2	ZX7	B	902	23/23	0.92	0.11	27,37,62,63	0

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.

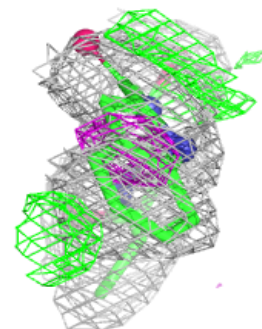
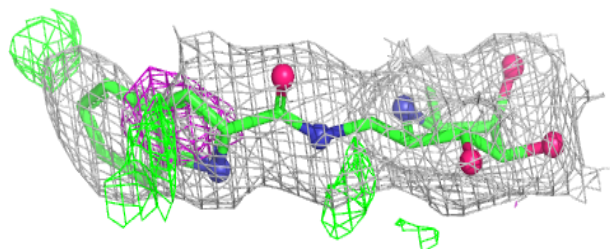
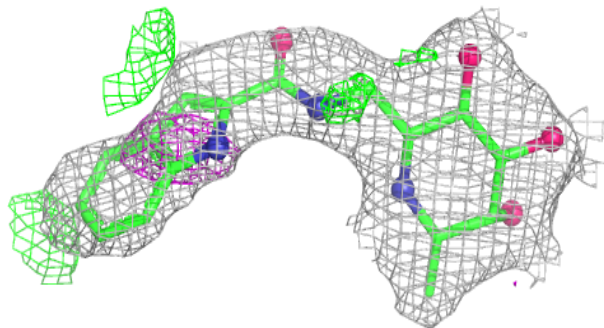
Electron density around ZX7 A 901:

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



Electron density around ZX7 B 902:

$2mF_o-DF_c$ (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.