



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 02:33 PM UTC

PDB ID : 6X3J / pdb_00006x3j
Title : Crystal structure of streptogramin A acetyltransferase VatA from Staphylococcus aureus in complex with streptogramin analog F0224 (46)
Authors : Chaires, H.A.; Fraser, J.S.
Deposited on : 2020-05-21
Resolution : 2.70 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)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

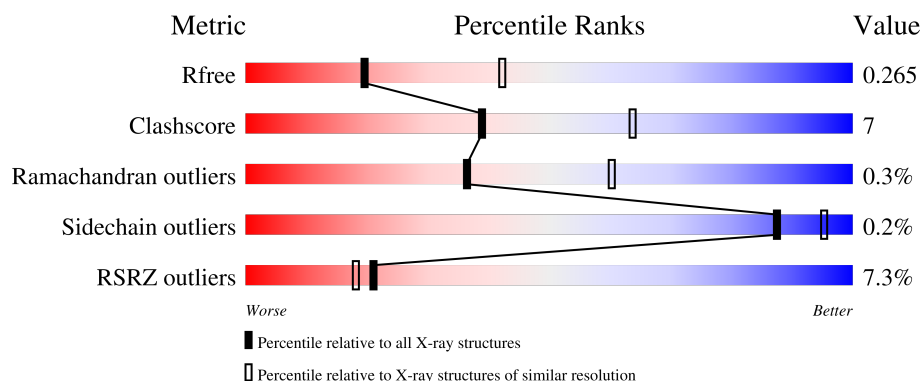
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	209	5% 85% 14% .
1	B	209	7% 80% 18% .
1	C	209	7% 89% 11% .
1	D	209	8% 85% 14% .
1	E	209	8% 84% 15% .

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Mol	Chain	Length	Quality of chain
1	F	209	

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
4	SXA	A	403	X	-	-	-
4	SXA	B	405	X	-	-	-
4	SXA	D	302	X	-	-	-
4	SXA	E	407	X	-	-	-
4	SXA	F	405	X	-	-	-
4	SXA	F	406	X	-	-	-

2 Entry composition [i](#)

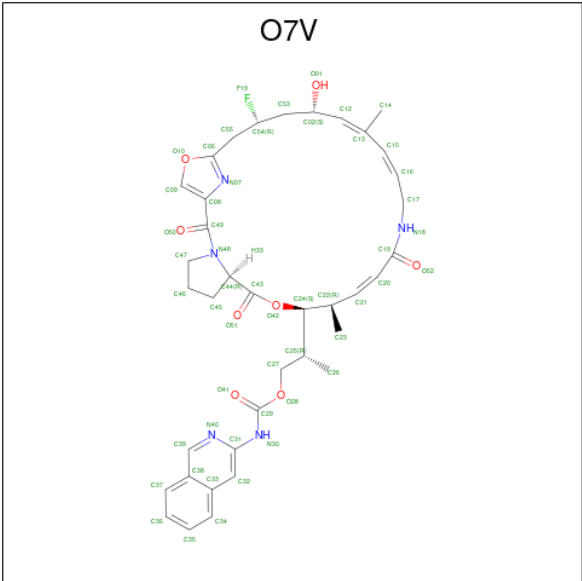
There are 7 unique types of molecules in this entry. The entry contains 20998 atoms, of which 10359 are hydrogens and 0 are deuteriums.

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 Virginiamycin A acetyltransferase,VatA.

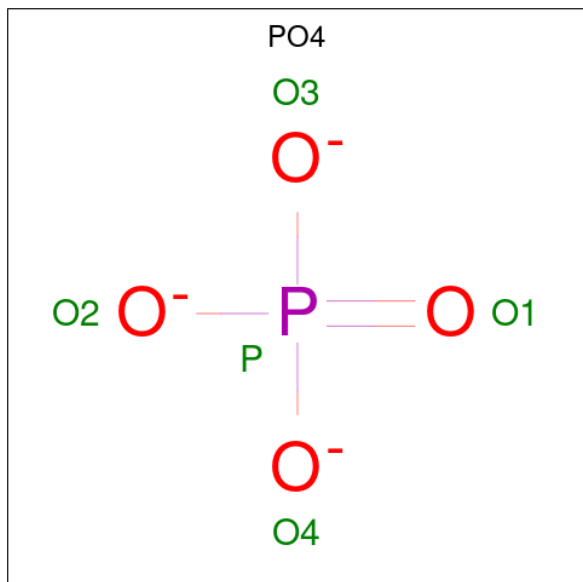
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	206	Total	C	H	N	O	S	0	0	0
			3295	1066	1646	277	298	8			
1	B	206	Total	C	H	N	O	S	0	0	0
			3295	1066	1646	277	298	8			
1	C	209	Total	C	H	N	O	S	0	0	0
			3359	1084	1682	284	301	8			
1	D	207	Total	C	H	N	O	S	0	0	0
			3315	1072	1655	281	299	8			
1	E	207	Total	C	H	N	O	S	0	0	0
			3319	1072	1659	281	299	8			
1	F	204	Total	C	H	N	O	S	0	0	0
			3248	1054	1619	271	296	8			

- Molecule 2 is (2R)-2-[(3S,4R,5E,10E,12E,14S,16R,26aR)-16-fluoro-14-hydroxy-4,12-dimethyl-1,7,22-trioxo-4,7,8,9,14,15,16,17,24,25,26,26a-dodecahydro-1H,3H,22H-21,18-(azeno)pyrrolo[2,1-c][1,8,4,19]dioxadiazacyclotetracosin-3-yl]propyl isoquinolin-3-ylcarbamate (CCD ID: O7V) (formula: C₃₈H₄₄FN₅O₈) (labeled as "Ligand of Interest" by depositor).



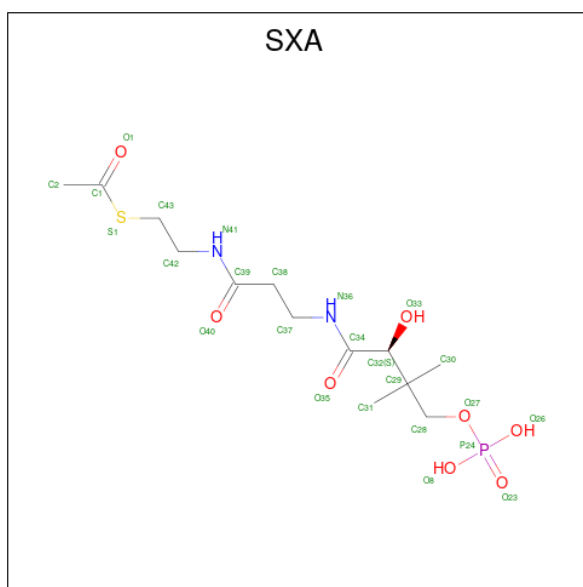
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	H	N	O	0	0
			96	38	1	44	5	8		
2	B	1	Total	C	F	H	N	O	2	0
			96	38	1	44	5	8		
2	B	1	Total	C	F	H	N	O	0	0
			96	38	1	44	5	8		
2	C	1	Total	C	F	H	N	O	0	1
			192	76	2	88	10	16		
2	E	1	Total	C	F	H	N	O	0	0
			96	38	1	44	5	8		
2	F	1	Total	C	F	H	N	O	0	0
			96	38	1	44	5	8		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

- Molecule 4 is THIOACETIC ACID S-{2-[3-(2-HYDROXY-3,3-DIMETHYL-4-PHOSPHO NOOXY-BUTYRYLAMINO)-PROPYONYLAMINO]-ETHYL} ESTER (CCD ID: SXA) (formula: C₁₃H₂₅N₂O₈PS).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total 44	C 13	H 24	N 2	O 4	S 1	0	0
4	B	1	Total 44	C 13	H 24	N 2	O 4	S 1	2	0
4	D	1	Total 44	C 13	H 24	N 2	O 4	S 1	2	0
4	E	1	Total 44	C 13	H 24	N 2	O 4	S 1	2	0
4	F	1	Total 44	C 13	H 24	N 2	O 4	S 1	2	0
4	F	1	Total 44	C 13	H 24	N 2	O 4	S 1	1	0

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total	Cl	0	0
			2	2		
5	C	2	Total	Cl	0	0
			2	2		
5	E	4	Total	Cl	0	0
			4	4		
5	F	3	Total	Cl	0	0
			3	3		

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	1	Total 1	Mg 1	0	0
6	D	1	Total 1	Mg 1	0	0

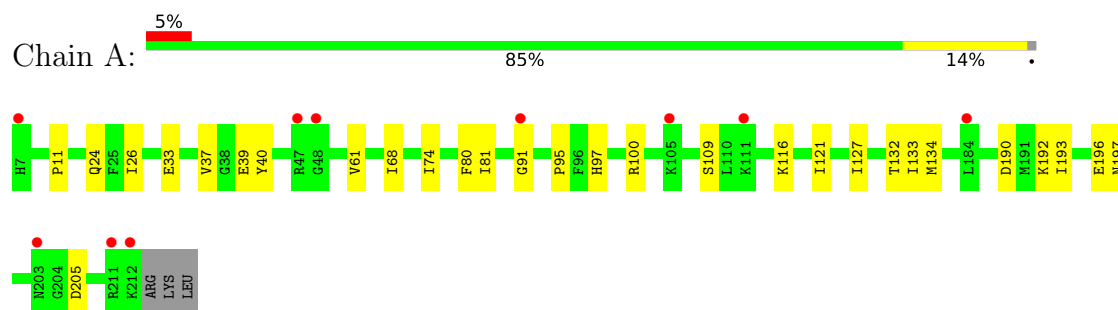
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	36	Total 36	O 36	0	0
7	B	45	Total 45	O 45	0	0
7	C	32	Total 32	O 32	0	0
7	D	28	Total 28	O 28	0	0
7	E	31	Total 31	O 31	0	0
7	F	36	Total 36	O 36	0	0

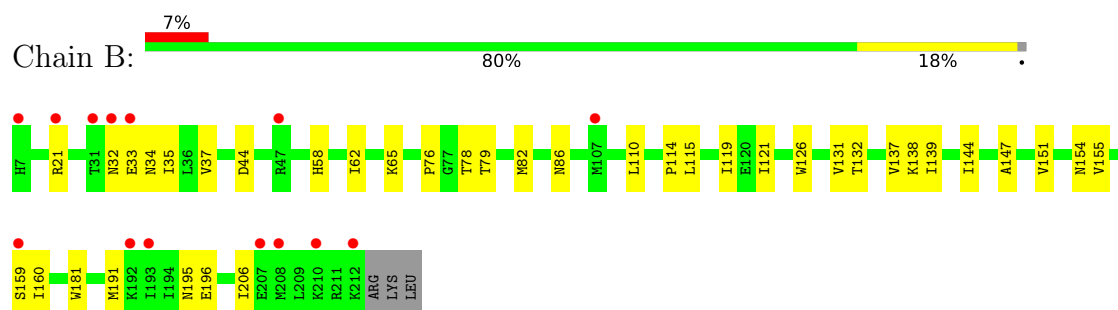
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

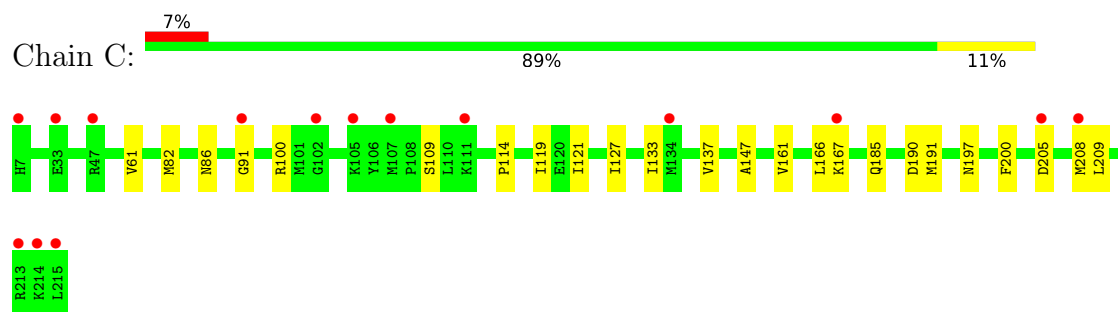
- Molecule 1: Virginiamycin A acetyltransferase,VatA



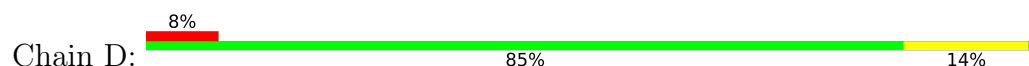
- Molecule 1: Virginiamycin A acetyltransferase,VatA

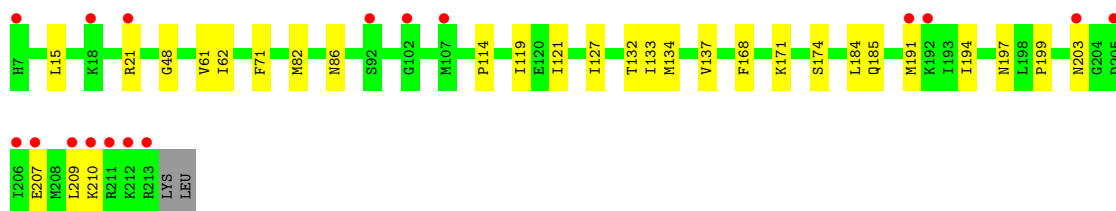


- Molecule 1: Virginiamycin A acetyltransferase,VatA

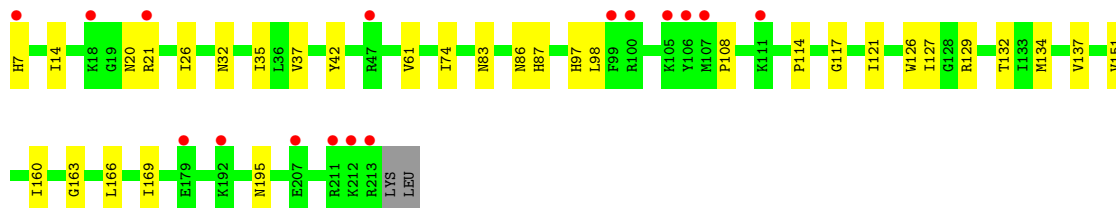
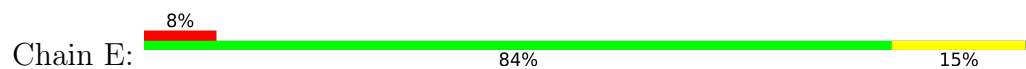


- Molecule 1: Virginiamycin A acetyltransferase,VatA

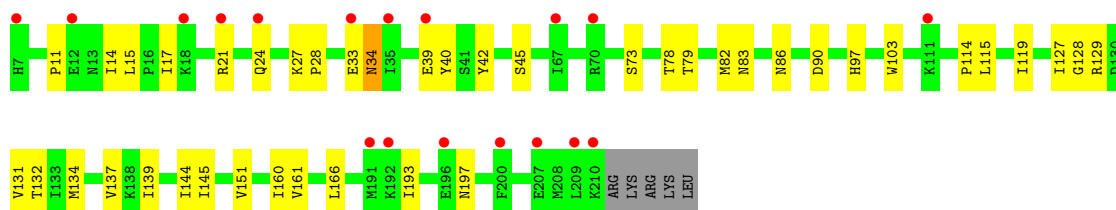
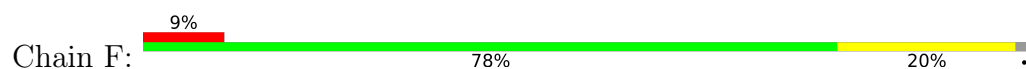




- Molecule 1: Virginiamycin A acetyltransferase, Vata



- Molecule 1: Virginiamycin A acetyltransferase, Vata



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	104.80Å 109.22Å 173.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	86.51 – 2.70 86.51 – 2.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (86.51-2.70) 100.0 (86.51-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.24 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.234 , 0.271 0.232 , 0.265	Depositor DCC
R_{free} test set	2797 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	53.4	Xtriage
Anisotropy	0.211	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.021 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	20998	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.94% 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: SXA, O7V, PO4, CL, MG

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	0.33	0/1691	0.53	1/2289 (0.0%)
1	B	0.44	0/1691	0.68	5/2289 (0.2%)
1	C	0.27	0/1719	0.49	1/2325 (0.0%)
1	D	0.31	0/1702	0.52	0/2303
1	E	0.31	0/1702	0.50	1/2303 (0.0%)
1	F	0.32	0/1671	0.58	4/2264 (0.2%)
All	All	0.33	0/10176	0.56	12/13773 (0.1%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	37	VAL	CA-C-N	6.27	127.62	120.53
1	B	37	VAL	C-N-CA	6.27	127.62	120.53
1	F	33	GLU	CA-C-O	-6.22	114.19	121.16
1	F	33	GLU	CA-C-N	6.13	133.25	121.54
1	F	33	GLU	C-N-CA	6.13	133.25	121.54
1	B	33	GLU	CA-C-N	5.62	132.28	121.54
1	B	33	GLU	C-N-CA	5.62	132.28	121.54
1	F	34	ASN	N-CA-CB	5.46	119.72	110.49
1	B	33	GLU	CA-C-O	-5.45	115.62	121.56
1	E	97	HIS	CA-CB-CG	-5.43	108.36	113.80
1	C	100	ARG	CB-CA-C	5.15	120.67	110.42
1	A	100	ARG	CB-CA-C	5.05	120.47	110.42

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	1649	1646	1645	16	5
1	B	1649	1646	1645	28	2
1	C	1677	1682	1682	14	5
1	D	1660	1655	1658	23	2
1	E	1660	1659	1658	27	0
1	F	1629	1619	1619	29	0
2	A	52	44	0	3	0
2	B	104	88	0	1	0
2	C	104	88	0	3	0
2	E	52	44	0	3	0
2	F	52	44	0	0	0
3	A	5	0	0	0	0
3	E	5	0	0	1	0
4	A	20	24	21	0	0
4	B	20	24	21	3	0
4	D	20	24	21	0	0
4	E	20	24	21	2	0
4	F	40	48	42	4	0
5	B	2	0	0	1	0
5	C	2	0	0	0	0
5	E	4	0	0	2	0
5	F	3	0	0	2	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
7	A	36	0	0	1	0
7	B	45	0	0	1	0
7	C	32	0	0	1	0
7	D	28	0	0	0	0
7	E	31	0	0	0	0
7	F	36	0	0	2	0
All	All	10639	10359	10033	139	7

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (139) 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:B:82:MET:HE1	1:B:119:ILE:HD12	1.32	1.06
1:B:82:MET:CE	1:B:119:ILE:HD12	1.87	1.04
1:E:87:HIS:NE2	4:F:405:SXA:O1	2.16	0.78
1:B:155:VAL:HG12	1:B:159:SER:OG	1.83	0.77
5:F:402:CL:CL	7:F:523:HOH:O	2.40	0.76
1:F:127:ILE:HG12	1:F:145:ILE:HD12	1.68	0.76
1:E:61:VAL:HG11	2:E:401:O7V:C26	2.18	0.74
1:A:81:ILE:HD12	1:A:134:MET:HE1	1.71	0.72
1:D:199:PRO:O	1:D:203:ASN:ND2	2.24	0.71
1:B:155:VAL:CG1	1:B:159:SER:OG	2.41	0.69
1:B:82:MET:HE2	1:B:119:ILE:HD12	1.75	0.69
5:B:403:CL:CL	7:B:528:HOH:O	2.49	0.67
1:E:132:THR:HG22	1:E:134:MET:CE	2.24	0.67
1:B:126:TRP:HE1	4:B:405:SXA:H2A	1.64	0.63
1:D:48:GLY:O	1:F:21:ARG:NH2	2.33	0.62
1:D:71:PHE:HZ	1:D:191:MET:HE1	1.64	0.62
1:D:21:ARG:HG2	1:D:21:ARG:O	2.00	0.61
1:F:193:ILE:O	1:F:197:ASN:ND2	2.35	0.60
1:F:79:THR:OG1	1:F:132:THR:HG23	2.02	0.60
1:F:21:ARG:O	1:F:21:ARG:HG2	2.02	0.59
1:F:86:ASN:HB3	1:F:114:PRO:HG2	1.85	0.59
1:A:11:PRO:HG3	1:A:40:TYR:CZ	2.39	0.58
1:D:48:GLY:O	1:F:21:ARG:CZ	2.52	0.58
1:B:110:LEU:HB3	1:B:115:LEU:HD11	1.85	0.57
1:F:83:ASN:HB3	1:F:115:LEU:CD2	2.34	0.57
1:C:197:ASN:HB3	1:C:209:LEU:HD21	1.87	0.56
1:E:86:ASN:HB3	1:E:114:PRO:HG2	1.86	0.56
1:F:83:ASN:HB3	1:F:115:LEU:HD22	1.88	0.56
1:F:14:ILE:HG22	1:F:15:LEU:HG	1.87	0.56
1:E:132:THR:HG22	1:E:134:MET:HE3	1.88	0.55
1:B:155:VAL:HG12	1:B:159:SER:HG	1.71	0.55
1:C:161:VAL:HG23	1:C:166:LEU:HD11	1.89	0.55
1:E:121:ILE:HD13	1:E:127:ILE:HD11	1.89	0.55
1:F:137:VAL:HG11	1:F:151:VAL:HG12	1.89	0.54
1:C:133:ILE:HG23	1:C:137:VAL:HG21	1.89	0.54
1:E:126:TRP:NE1	4:E:407:SXA:S1	2.75	0.54
1:D:121:ILE:HD13	1:D:127:ILE:HD11	1.88	0.54
1:D:207:GLU:HA	1:D:210:LYS:HB2	1.89	0.54
1:B:86:ASN:HB3	1:B:114:PRO:HG2	1.90	0.54
1:E:98:LEU:HB3	1:F:17:ILE:HD11	1.90	0.53
1:D:133:ILE:HG23	1:D:137:VAL:HG21	1.90	0.53
1:E:32:ASN:HB3	1:E:35:ILE:HD12	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:163:GLY:N	5:E:405:CL:CL	2.77	0.52
1:B:126:TRP:NE1	4:B:405:SXA:H2A	2.25	0.52
1:E:108:PRO:CG	2:E:401:O7V:O52	2.57	0.52
1:E:26:ILE:HB	1:E:37:VAL:HG11	1.92	0.52
1:F:24:GLN:NE2	1:F:45:SER:HB3	2.25	0.52
1:B:44:ASP:O	1:B:76:PRO:HB3	2.10	0.51
1:C:185:GLN:NE2	7:C:501:HOH:O	2.40	0.51
1:D:61:VAL:HG23	1:D:62:ILE:HG13	1.91	0.51
1:D:15:LEU:HD22	1:D:21:ARG:HA	1.93	0.51
1:E:137:VAL:HG11	1:E:151:VAL:HG12	1.92	0.51
1:E:74:ILE:CD1	1:E:127:ILE:HD12	2.41	0.50
1:D:132:THR:HG22	1:D:134:MET:HE2	1.93	0.49
1:A:97:HIS:NE2	2:A:401:O7V:O52	2.45	0.49
1:F:132:THR:HG22	1:F:134:MET:HE2	1.95	0.49
1:D:184:LEU:HD21	1:D:209:LEU:HD13	1.94	0.48
1:E:151:VAL:HG13	1:E:166:LEU:HD22	1.95	0.48
1:D:82:MET:SD	1:D:119:ILE:HD12	2.53	0.48
1:F:82:MET:SD	1:F:119:ILE:HD12	2.53	0.48
1:B:44:ASP:HB3	1:B:76:PRO:HD3	1.96	0.48
1:F:97:HIS:HA	1:F:103:TRP:O	2.14	0.48
1:B:78:THR:HA	1:B:131:VAL:O	2.14	0.47
1:B:147:ALA:N	4:B:405:SXA:O40	2.43	0.47
1:B:34:ASN:O	1:B:65:LYS:HA	2.14	0.47
1:E:137:VAL:HG11	1:E:151:VAL:CG1	2.44	0.47
1:E:98:LEU:HB3	1:F:17:ILE:CD1	2.45	0.47
1:A:61:VAL:HG11	2:A:401:O7V:C31	2.44	0.47
1:C:167:LYS:HE3	7:F:522:HOH:O	2.14	0.47
1:B:21:ARG:HH11	1:E:21:ARG:HH22	1.62	0.47
1:A:121:ILE:HD13	1:A:127:ILE:HD11	1.96	0.46
1:C:147:ALA:HB2	4:F:406:SXA:C42	2.45	0.46
1:B:121:ILE:HG12	1:B:139:ILE:HD12	1.98	0.46
1:B:181:TRP:CD2	1:B:206:ILE:HD13	2.51	0.46
1:C:61:VAL:HG22	1:C:61:VAL:O	2.16	0.46
1:D:191:MET:O	1:D:191:MET:HG3	2.15	0.46
1:D:197:ASN:HB3	1:D:209:LEU:HD21	1.96	0.46
1:C:190:ASP:OD2	1:C:191:MET:N	2.49	0.46
1:A:190:ASP:OD1	1:A:192:LYS:HD3	2.16	0.45
1:A:39:GLU:O	1:A:40:TYR:HB2	2.16	0.45
1:D:199:PRO:O	1:D:203:ASN:CG	2.59	0.45
2:A:401:O7V:C47	2:A:401:O7V:N07	2.79	0.45
1:C:121:ILE:HD13	1:C:127:ILE:HD11	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:14:ILE:N	3:E:406:PO4:O4	2.29	0.45
1:F:90:ASP:N	5:F:403:CL:CL	2.87	0.45
1:A:132:THR:HG22	1:A:134:MET:CE	2.47	0.45
1:F:27:LYS:HD2	1:F:39:GLU:HG2	1.98	0.44
1:E:83:ASN:ND2	1:E:117:GLY:O	2.50	0.44
1:B:34:ASN:HB3	1:B:65:LYS:HG2	2.00	0.44
1:B:191:MET:O	1:B:195:ASN:ND2	2.50	0.43
1:D:133:ILE:HG23	1:D:137:VAL:CG2	2.48	0.43
1:B:79:THR:OG1	1:B:132:THR:HG23	2.18	0.43
2:C:401[A]:O7V:C47	2:C:401[A]:O7V:N07	2.81	0.43
1:D:86:ASN:HB3	1:D:114:PRO:HG2	2.00	0.43
1:D:132:THR:CG2	1:D:134:MET:HE2	2.48	0.43
1:C:82:MET:SD	1:C:119:ILE:HD12	2.59	0.43
1:C:133:ILE:HG23	1:C:137:VAL:CG2	2.48	0.43
1:B:35:ILE:HG13	1:B:58:HIS:HE1	1.83	0.43
1:F:78:THR:HA	1:F:131:VAL:O	2.19	0.43
1:F:128:GLY:HA2	4:F:405:SXA:C42	2.49	0.42
1:F:144:ILE:HB	1:F:160:ILE:HG12	2.00	0.42
1:A:192:LYS:N	1:A:192:LYS:HD2	2.34	0.42
1:F:39:GLU:O	1:F:40:TYR:HB2	2.20	0.42
1:C:161:VAL:CG2	1:C:166:LEU:HD11	2.49	0.42
1:F:129:ARG:H	4:F:405:SXA:H42A	1.84	0.42
1:A:24:GLN:NE2	7:A:502:HOH:O	2.52	0.42
2:C:401[A]:O7V:O01	1:E:42:TYR:OH	2.31	0.42
1:F:11:PRO:HB2	1:F:39:GLU:HB3	2.02	0.42
1:A:26:ILE:HB	1:A:37:VAL:HG11	2.02	0.42
1:A:33:GLU:H	1:A:33:GLU:CD	2.28	0.42
1:B:137:VAL:HG11	1:B:151:VAL:HG12	2.02	0.42
1:C:200:PHE:CE1	1:C:208:MET:HE1	2.55	0.42
1:F:42:TYR:CZ	1:F:73:SER:HB2	2.55	0.42
2:B:402:O7V:N07	2:B:402:O7V:C47	2.83	0.41
1:B:191:MET:HE2	1:B:191:MET:HA	2.01	0.41
1:E:7:HIS:O	1:E:195:ASN:ND2	2.51	0.41
1:E:129:ARG:H	4:E:407:SXA:H42	1.84	0.41
1:A:80:PHE:CE1	1:A:133:ILE:HD12	2.55	0.41
1:A:68:ILE:HD13	1:A:74:ILE:HD11	2.03	0.41
1:C:86:ASN:HB3	1:C:114:PRO:HG2	2.03	0.41
1:B:62:ILE:HG12	1:B:110:LEU:HD21	2.01	0.41
1:B:144:ILE:HB	1:B:160:ILE:HG12	2.03	0.41
1:D:184:LEU:C	1:D:185:GLN:HG2	2.46	0.41
1:E:20:ASN:O	5:E:404:CL:CL	2.76	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:21:ARG:O	1:E:21:ARG:HG3	2.21	0.41
1:F:139:ILE:HD13	1:F:145:ILE:HD11	2.03	0.41
1:B:126:TRP:CE3	1:B:144:ILE:HG23	2.56	0.41
2:E:401:O7V:N07	2:E:401:O7V:C47	2.84	0.41
1:F:27:LYS:N	1:F:28:PRO:HD2	2.36	0.41
1:D:168:PHE:HZ	1:D:171:LYS:HG3	1.86	0.40
1:E:160:ILE:HG22	1:E:169:ILE:HD12	2.03	0.40
1:B:138:LYS:H	1:B:154:ASN:ND2	2.19	0.40
1:A:192:LYS:O	1:A:196:GLU:HG3	2.21	0.40
1:F:161:VAL:HG23	1:F:166:LEU:HD11	2.04	0.40
1:A:193:ILE:O	1:A:197:ASN:ND2	2.55	0.40
1:E:26:ILE:HG21	1:E:37:VAL:HG21	2.03	0.40
2:C:401[B]:O7V:N07	2:C:401[B]:O7V:C47	2.85	0.40
1:D:71:PHE:CZ	1:D:191:MET:HE1	2.50	0.40
1:D:191:MET:SD	1:D:194:ILE:HB	2.62	0.40

All (7) 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)
1:A:109:SER:OG	1:C:205:ASP:OD1[4_545]	1.03	1.17
1:B:196:GLU:OE1	1:D:174:SER:HG[3_644]	1.05	0.55
1:A:109:SER:HG	1:C:205:ASP:OD1[4_545]	1.11	0.49
1:B:196:GLU:OE1	1:D:174:SER:OG[3_644]	1.86	0.34
1:A:109:SER:OG	1:C:205:ASP:CG[4_545]	2.01	0.19
1:A:205:ASP:OD2	1:C:109:SER:OG[3_554]	2.01	0.19
1:A:109:SER:HG	1:C:205:ASP:CG[4_545]	1.58	0.02

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	204/209 (98%)	197 (97%)	5 (2%)	2 (1%)	12	32
1	B	204/209 (98%)	201 (98%)	3 (2%)	0	100	100
1	C	207/209 (99%)	199 (96%)	7 (3%)	1 (0%)	24	48
1	D	205/209 (98%)	199 (97%)	6 (3%)	0	100	100
1	E	205/209 (98%)	200 (98%)	5 (2%)	0	100	100
1	F	202/209 (97%)	197 (98%)	4 (2%)	1 (0%)	24	48
All	All	1227/1254 (98%)	1193 (97%)	30 (2%)	4 (0%)	36	60

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	34	ASN
1	A	91	GLY
1	C	91	GLY
1	A	95	PRO

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	178/181 (98%)	177 (99%)	1 (1%)	78	91
1	B	178/181 (98%)	177 (99%)	1 (1%)	78	91
1	C	181/181 (100%)	181 (100%)	0	100	100
1	D	179/181 (99%)	179 (100%)	0	100	100
1	E	179/181 (99%)	179 (100%)	0	100	100
1	F	176/181 (97%)	176 (100%)	0	100	100
All	All	1071/1086 (99%)	1069 (100%)	2 (0%)	87	95

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	116	LYS

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Mol	Chain	Res	Type
1	B	32	ASN

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

Mol	Chain	Res	Type
1	B	24	GLN
1	B	58	HIS
1	B	154	ASN
1	C	24	GLN
1	C	54	GLN
1	C	164	ASN
1	F	24	GLN

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](#)

Of 28 ligands modelled in this entry, 13 are monoatomic - leaving 15 for Mogul analysis.

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	O7V	C	401[B]	-	54,56,56	1.29	4 (7%)	67,77,77	1.33	10 (14%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	O7V	B	402	-	54,56,56	1.30	4 (7%)	67,77,77	1.16	9 (13%)
3	PO4	A	402	-	4,4,4	0.95	0	6,6,6	0.50	0
4	SXA	D	302	-	18,19,24	1.84	1 (5%)	23,25,33	1.28	3 (13%)
4	SXA	F	405	-	18,19,24	1.75	1 (5%)	23,25,33	1.24	1 (4%)
2	O7V	E	401	-	54,56,56	1.35	5 (9%)	67,77,77	1.69	14 (20%)
3	PO4	E	406	-	4,4,4	0.97	0	6,6,6	0.46	0
4	SXA	A	403	-	18,19,24	1.43	1 (5%)	23,25,33	0.76	0
4	SXA	F	406	-	18,19,24	1.49	1 (5%)	23,25,33	1.04	2 (8%)
4	SXA	B	405	-	18,19,24	1.42	1 (5%)	23,25,33	1.21	3 (13%)
2	O7V	F	401	-	54,56,56	1.32	5 (9%)	67,77,77	1.20	7 (10%)
2	O7V	A	401	-	54,56,56	1.35	5 (9%)	67,77,77	1.36	8 (11%)
2	O7V	C	401[A]	-	54,56,56	1.29	3 (5%)	67,77,77	1.35	10 (14%)
4	SXA	E	407	-	18,19,24	1.37	1 (5%)	23,25,33	1.04	1 (4%)
2	O7V	B	401	-	54,56,56	1.33	5 (9%)	67,77,77	1.24	8 (11%)

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	O7V	C	401[B]	-	-	15/57/67/67	0/4/5/5
2	O7V	B	402	-	-	10/57/67/67	0/4/5/5
4	SXA	D	302	-	1/1/7/8	13/23/23/30	-
2	O7V	E	401	-	-	13/57/67/67	0/4/5/5
2	O7V	A	401	-	-	10/57/67/67	0/4/5/5
4	SXA	A	403	-	1/1/7/8	6/23/23/30	-
4	SXA	F	406	-	1/1/7/8	5/23/23/30	-
4	SXA	B	405	-	1/1/7/8	8/23/23/30	-
2	O7V	F	401	-	-	10/57/67/67	0/4/5/5
4	SXA	F	405	-	1/1/7/8	11/23/23/30	-
2	O7V	C	401[A]	-	-	11/57/67/67	0/4/5/5
4	SXA	E	407	-	1/1/7/8	8/23/23/30	-
2	O7V	B	401	-	-	11/57/67/67	0/4/5/5

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	302	SXA	C29-C32	6.96	1.59	1.55
4	F	405	SXA	C29-C32	6.64	1.59	1.55
4	F	406	SXA	C29-C32	5.32	1.58	1.55
4	A	403	SXA	C29-C32	5.08	1.58	1.55
4	B	405	SXA	C29-C32	4.97	1.58	1.55
4	E	407	SXA	C29-C32	4.92	1.58	1.55
2	C	401[B]	O7V	C09-C08	4.86	1.39	1.34
2	B	402	O7V	C09-C08	4.83	1.39	1.34
2	C	401[A]	O7V	C09-C08	4.82	1.39	1.34
2	F	401	O7V	C09-C08	4.81	1.39	1.34
2	A	401	O7V	C09-C08	4.79	1.39	1.34
2	B	401	O7V	C09-C08	4.75	1.39	1.34
2	E	401	O7V	C09-C08	4.49	1.39	1.34
2	E	401	O7V	C08-C49	-3.37	1.44	1.49
2	A	401	O7V	C08-N07	-3.19	1.31	1.38
2	C	401[A]	O7V	C08-N07	-3.09	1.32	1.38
2	E	401	O7V	C08-N07	-3.08	1.32	1.38
2	A	401	O7V	C08-C49	-2.87	1.45	1.49
2	B	401	O7V	C08-N07	-2.87	1.32	1.38
2	B	402	O7V	C08-N07	-2.86	1.32	1.38
2	C	401[A]	O7V	C08-C49	-2.77	1.45	1.49
2	B	401	O7V	C08-C49	-2.74	1.45	1.49
2	C	401[B]	O7V	C08-N07	-2.72	1.32	1.38
2	F	401	O7V	C08-N07	-2.59	1.33	1.38
2	B	402	O7V	C08-C49	-2.50	1.45	1.49
2	C	401[B]	O7V	C20-C19	-2.47	1.43	1.48
2	A	401	O7V	C02-C12	2.44	1.53	1.50
2	C	401[B]	O7V	C08-C49	-2.40	1.46	1.49
2	F	401	O7V	C06-N07	2.32	1.31	1.29
2	F	401	O7V	C08-C49	-2.25	1.46	1.49
2	E	401	O7V	C02-C12	2.15	1.53	1.50
2	F	401	O7V	C39-N40	2.13	1.36	1.32
2	E	401	O7V	O42-C24	-2.12	1.41	1.44
2	B	402	O7V	C39-N40	2.12	1.36	1.32
2	A	401	O7V	C39-N40	2.08	1.36	1.32
2	B	401	O7V	C39-N40	2.03	1.35	1.32
2	B	401	O7V	C20-C19	-2.03	1.44	1.48

All (76) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	401	O7V	C17-N18-C19	6.55	132.79	122.17
2	A	401	O7V	C27-O28-C29	5.07	128.04	115.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401[B]	O7V	O28-C27-C25	4.66	115.95	108.28
2	A	401	O7V	C53-C02-C12	4.24	117.49	111.28
2	E	401	O7V	O42-C24-C25	-4.22	102.31	108.15
2	C	401[A]	O7V	C27-O28-C29	4.17	125.88	115.91
2	C	401[A]	O7V	C53-C02-C12	4.10	117.29	111.28
2	E	401	O7V	O28-C27-C25	3.90	114.69	108.28
2	E	401	O7V	O52-C19-N18	-3.82	116.29	122.29
2	B	401	O7V	C27-O28-C29	3.79	124.96	115.91
2	F	401	O7V	C27-O28-C29	3.77	124.93	115.91
2	C	401[A]	O7V	O28-C27-C25	3.49	114.03	108.28
2	B	402	O7V	C27-O28-C29	3.46	124.18	115.91
2	E	401	O7V	C53-C02-C12	3.31	116.13	111.28
2	A	401	O7V	O28-C27-C25	3.30	113.70	108.28
2	E	401	O7V	C24-O42-C43	3.29	123.23	117.76
2	A	401	O7V	O28-C29-N30	3.23	114.42	109.35
2	C	401[B]	O7V	C27-C25-C24	3.21	116.25	111.08
2	B	401	O7V	O28-C27-C25	3.13	113.42	108.28
4	F	405	SXA	C43-C42-N41	3.12	118.94	112.41
2	E	401	O7V	C23-C22-C24	3.01	116.41	111.11
4	B	405	SXA	C38-C37-N36	3.00	118.39	112.00
2	C	401[B]	O7V	C27-O28-C29	2.93	122.91	115.91
2	B	401	O7V	C53-C02-C12	2.88	115.51	111.28
4	D	302	SXA	C43-C42-N41	2.81	118.29	112.41
2	F	401	O7V	C49-C08-N07	2.73	128.48	123.70
2	E	401	O7V	C27-O28-C29	2.69	122.34	115.91
2	C	401[A]	O7V	O28-C29-N30	2.69	113.57	109.35
2	B	401	O7V	O28-C29-N30	2.66	113.53	109.35
4	F	406	SXA	C37-C38-C39	2.61	116.73	112.39
2	C	401[B]	O7V	C53-C02-C12	2.55	115.02	111.28
2	A	401	O7V	O28-C29-O41	-2.50	119.47	124.26
4	D	302	SXA	C38-C37-N36	2.49	117.31	112.00
2	E	401	O7V	C27-C25-C24	2.43	115.00	111.08
2	C	401[A]	O7V	C46-C47-N48	2.43	107.41	103.24
2	B	402	O7V	C53-C02-C12	2.42	114.82	111.28
4	D	302	SXA	C37-C38-C39	2.41	116.41	112.39
2	B	401	O7V	C16-C15-C13	-2.40	122.27	125.89
4	B	405	SXA	C43-C42-N41	2.39	117.40	112.41
2	F	401	O7V	O28-C29-N30	2.37	113.07	109.35
2	C	401[B]	O7V	C49-C08-N07	2.33	127.79	123.70
2	C	401[B]	O7V	O42-C24-C25	2.33	111.39	108.15
2	B	401	O7V	C49-C08-N07	2.33	127.78	123.70
2	C	401[B]	O7V	C46-C47-N48	2.31	107.21	103.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	O7V	C46-C47-N48	2.29	107.17	103.24
2	C	401[B]	O7V	C23-C22-C24	2.26	115.09	111.11
2	C	401[A]	O7V	C24-O42-C43	2.26	121.52	117.76
2	F	401	O7V	O28-C29-O41	-2.25	119.94	124.26
2	B	401	O7V	C24-O42-C43	2.25	121.51	117.76
2	B	402	O7V	C24-O42-C43	2.25	121.50	117.76
2	B	402	O7V	C23-C22-C21	2.23	114.95	109.91
2	B	402	O7V	O28-C29-N30	2.21	112.83	109.35
4	B	405	SXA	C37-N36-C34	2.18	126.46	122.55
2	B	402	O7V	C26-C25-C24	2.17	114.69	111.14
2	B	402	O7V	C46-C47-N48	2.17	106.97	103.24
2	A	401	O7V	C09-C08-N07	-2.17	107.54	109.15
2	E	401	O7V	C47-N48-C44	-2.14	108.65	112.01
2	E	401	O7V	C26-C25-C27	2.14	113.94	110.03
4	F	406	SXA	C38-C37-N36	2.13	116.54	112.00
2	F	401	O7V	C09-C08-N07	-2.13	107.56	109.15
2	E	401	O7V	C16-C15-C13	-2.13	122.67	125.89
2	E	401	O7V	C20-C19-N18	2.13	119.73	115.07
2	C	401[B]	O7V	C31-N30-C29	2.13	133.07	127.42
2	A	401	O7V	C24-O42-C43	2.12	121.28	117.76
2	F	401	O7V	C23-C22-C21	2.10	114.66	109.91
2	C	401[A]	O7V	O28-C29-O41	-2.10	120.24	124.26
2	B	402	O7V	O28-C27-C25	2.08	111.70	108.28
4	E	407	SXA	C38-C37-N36	2.06	116.39	112.00
2	E	401	O7V	C45-C44-C43	2.04	115.76	111.21
2	C	401[A]	O7V	C14-C13-C15	-2.04	114.97	118.09
2	C	401[A]	O7V	C26-C25-C27	2.04	113.76	110.03
2	B	402	O7V	C09-C08-N07	-2.04	107.63	109.15
2	C	401[A]	O7V	C09-C08-N07	-2.02	107.64	109.15
2	F	401	O7V	C24-O42-C43	2.02	121.11	117.76
2	B	401	O7V	C17-N18-C19	2.02	125.44	122.17
2	C	401[B]	O7V	C09-C08-N07	-2.00	107.66	109.15

All (6) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	A	403	SXA	C32
4	B	405	SXA	C32
4	D	302	SXA	C32
4	E	407	SXA	C32
4	F	405	SXA	C32
4	F	406	SXA	C32

All (131) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	O7V	C26-C25-C27-O28
2	B	401	O7V	C12-C13-C15-C16
2	B	401	O7V	C14-C13-C15-C16
2	B	401	O7V	O42-C24-C25-C26
2	B	402	O7V	O42-C24-C25-C26
2	B	402	O7V	N30-C29-O28-C27
2	B	402	O7V	O41-C29-O28-C27
2	C	401[A]	O7V	C12-C13-C15-C16
2	C	401[A]	O7V	C14-C13-C15-C16
2	C	401[A]	O7V	C24-C25-C27-O28
2	C	401[A]	O7V	C26-C25-C27-O28
2	C	401[A]	O7V	N30-C29-O28-C27
2	C	401[A]	O7V	O41-C29-O28-C27
2	C	401[A]	O7V	F19-C54-C55-C06
2	C	401[B]	O7V	C12-C13-C15-C16
2	C	401[B]	O7V	C24-C25-C27-O28
2	C	401[B]	O7V	C26-C25-C27-O28
2	C	401[B]	O7V	N30-C29-O28-C27
2	C	401[B]	O7V	O41-C29-O28-C27
2	C	401[B]	O7V	F19-C54-C55-C06
2	E	401	O7V	C21-C22-C24-C25
2	E	401	O7V	C23-C22-C24-C25
2	E	401	O7V	C20-C19-N18-C17
2	E	401	O7V	O52-C19-N18-C17
2	E	401	O7V	C24-C25-C27-O28
2	E	401	O7V	C26-C25-C27-O28
2	F	401	O7V	O42-C24-C25-C26
2	F	401	O7V	N30-C29-O28-C27
2	F	401	O7V	O41-C29-O28-C27
4	A	403	SXA	N36-C37-C38-C39
4	A	403	SXA	N41-C42-C43-S1
4	B	405	SXA	N36-C37-C38-C39
4	B	405	SXA	O1-C1-S1-C43
4	B	405	SXA	C2-C1-S1-C43
4	D	302	SXA	C28-C29-C32-O33
4	D	302	SXA	C28-C29-C32-C34
4	D	302	SXA	C30-C29-C32-O33
4	D	302	SXA	C30-C29-C32-C34
4	D	302	SXA	C31-C29-C32-O33
4	D	302	SXA	C31-C29-C32-C34
4	D	302	SXA	N36-C37-C38-C39
4	D	302	SXA	C42-C43-S1-C1

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Mol	Chain	Res	Type	Atoms
4	D	302	SXA	O1-C1-S1-C43
4	D	302	SXA	C2-C1-S1-C43
4	E	407	SXA	O33-C32-C34-O35
4	E	407	SXA	N36-C37-C38-C39
4	E	407	SXA	C42-C43-S1-C1
4	E	407	SXA	O1-C1-S1-C43
4	E	407	SXA	C2-C1-S1-C43
4	F	405	SXA	C28-C29-C32-O33
4	F	405	SXA	C28-C29-C32-C34
4	F	405	SXA	C30-C29-C32-O33
4	F	405	SXA	C30-C29-C32-C34
4	F	405	SXA	C31-C29-C32-O33
4	F	405	SXA	C31-C29-C32-C34
4	F	405	SXA	N36-C37-C38-C39
4	F	405	SXA	O1-C1-S1-C43
4	F	405	SXA	C2-C1-S1-C43
4	F	406	SXA	N36-C37-C38-C39
4	F	406	SXA	N41-C42-C43-S1
2	A	401	O7V	O41-C29-O28-C27
2	B	401	O7V	O41-C29-O28-C27
2	A	401	O7V	N30-C29-O28-C27
2	B	401	O7V	N30-C29-O28-C27
2	E	401	O7V	O42-C43-C44-C45
2	C	401[B]	O7V	O52-C19-C20-C21
2	C	401[B]	O7V	N18-C19-C20-C21
2	B	402	O7V	C14-C13-C15-C16
2	C	401[B]	O7V	C14-C13-C15-C16
2	E	401	O7V	C14-C13-C15-C16
2	F	401	O7V	C14-C13-C15-C16
2	A	401	O7V	C20-C21-C22-C23
2	E	401	O7V	O51-C43-C44-C45
2	A	401	O7V	C12-C13-C15-C16
2	C	401[B]	O7V	C20-C21-C22-C23
4	A	403	SXA	O1-C1-S1-C43
4	A	403	SXA	C2-C1-S1-C43
2	C	401[B]	O7V	C20-C21-C22-C24
2	A	401	O7V	C14-C13-C15-C16
2	B	401	O7V	N18-C19-C20-C21
4	A	403	SXA	O33-C32-C34-O35
4	F	405	SXA	O33-C32-C34-O35
4	F	406	SXA	O33-C32-C34-O35
2	C	401[A]	O7V	O42-C24-C25-C26

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Mol	Chain	Res	Type	Atoms
2	A	401	O7V	O52-C19-C20-C21
2	C	401[A]	O7V	C53-C54-C55-C06
4	D	302	SXA	C29-C32-C34-O35
2	A	401	O7V	N18-C19-C20-C21
4	B	405	SXA	C29-C32-C34-N36
4	D	302	SXA	C29-C32-C34-N36
2	B	401	O7V	O52-C19-C20-C21
2	B	402	O7V	C12-C13-C15-C16
4	E	407	SXA	O33-C32-C34-N36
4	F	405	SXA	O33-C32-C34-N36
4	F	406	SXA	O33-C32-C34-N36
2	C	401[A]	O7V	C20-C21-C22-C23
2	B	402	O7V	C15-C16-C17-N18
4	B	405	SXA	O33-C32-C34-O35
2	E	401	O7V	C12-C13-C15-C16
2	E	401	O7V	C21-C22-C24-O42
2	E	401	O7V	C23-C22-C24-O42
4	B	405	SXA	C42-C43-S1-C1
4	F	406	SXA	C42-C43-S1-C1
2	B	401	O7V	C22-C24-C25-C26
2	B	401	O7V	C22-C24-C25-C27
2	B	402	O7V	C22-C24-C25-C26
2	F	401	O7V	C22-C24-C25-C26
2	A	401	O7V	C53-C54-C55-C06
2	C	401[B]	O7V	C53-C54-C55-C06
4	E	407	SXA	C29-C32-C34-O35
4	E	407	SXA	C29-C32-C34-N36
2	B	402	O7V	C20-C21-C22-C23
2	A	401	O7V	F19-C54-C55-C06
2	B	401	O7V	C15-C16-C17-N18
2	F	401	O7V	C15-C16-C17-N18
4	A	403	SXA	O33-C32-C34-N36
4	D	302	SXA	O33-C32-C34-N36
2	F	401	O7V	C20-C21-C22-C23
2	C	401[A]	O7V	O51-C43-C44-C45
2	B	402	O7V	C22-C24-C25-C27
2	F	401	O7V	C22-C24-C25-C27
2	B	401	O7V	C20-C21-C22-C23
2	C	401[B]	O7V	C15-C16-C17-N18
4	B	405	SXA	C29-C32-C34-O35
2	C	401[B]	O7V	O42-C43-C44-C45
2	F	401	O7V	C12-C13-C15-C16

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Mol	Chain	Res	Type	Atoms
2	C	401[B]	O7V	O10-C06-C55-C54
2	E	401	O7V	O42-C24-C25-C27
2	F	401	O7V	O10-C06-C55-C54
4	B	405	SXA	O33-C32-C34-N36
2	B	402	O7V	C22-C24-O42-C43

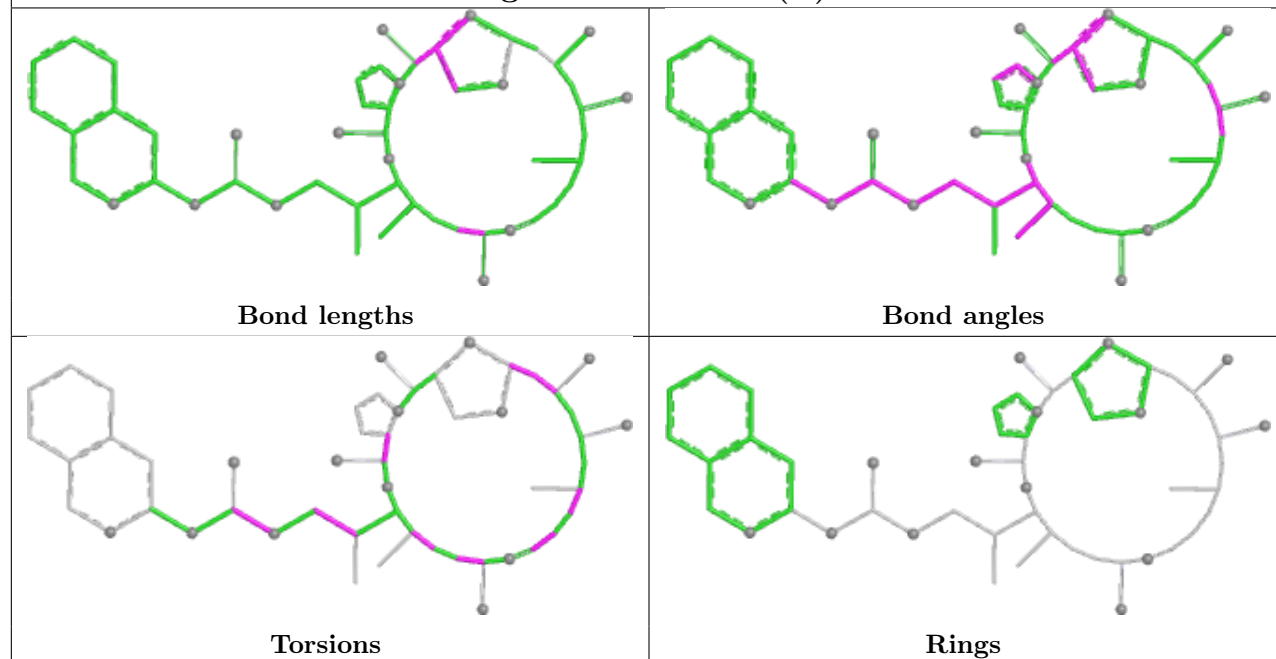
There are no ring outliers.

10 monomers are involved in 20 short contacts:

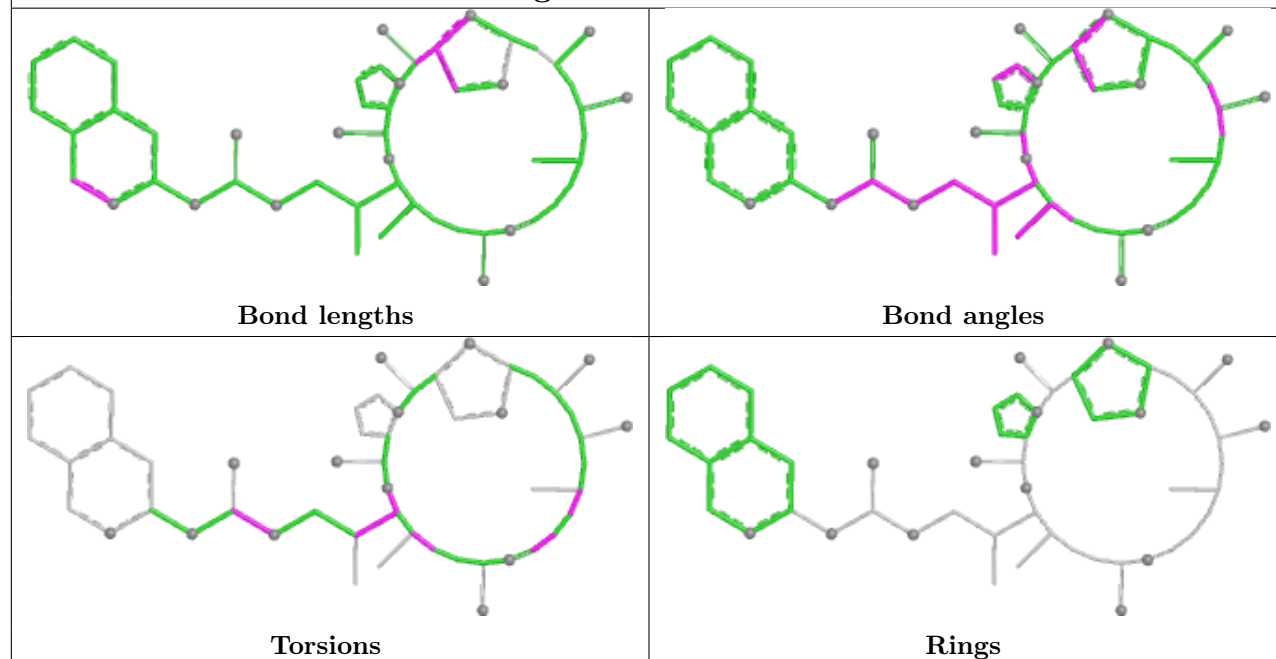
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401[B]	O7V	1	0
2	B	402	O7V	1	0
4	F	405	SXA	3	0
2	E	401	O7V	3	0
3	E	406	PO4	1	0
4	F	406	SXA	1	0
4	B	405	SXA	3	0
2	A	401	O7V	3	0
2	C	401[A]	O7V	2	0
4	E	407	SXA	2	0

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.

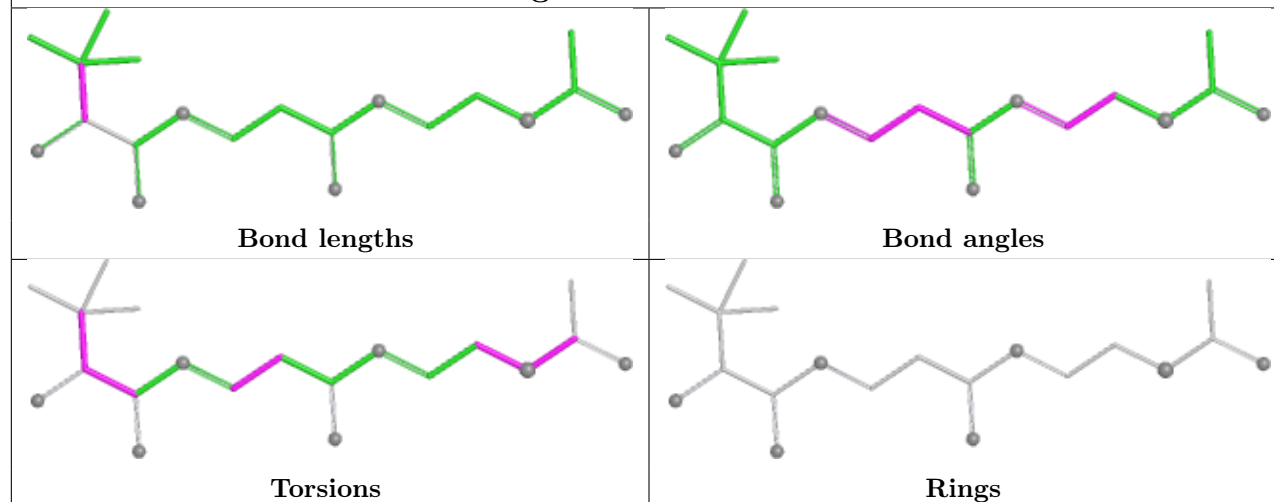
Ligand O7V C 401 (B)



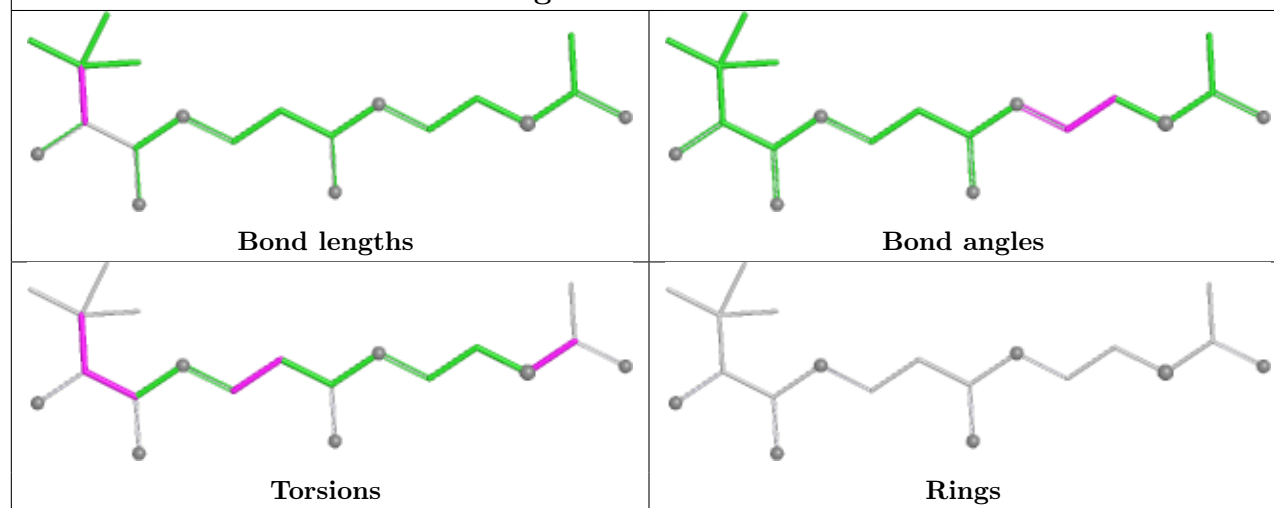
Ligand O7V B 402



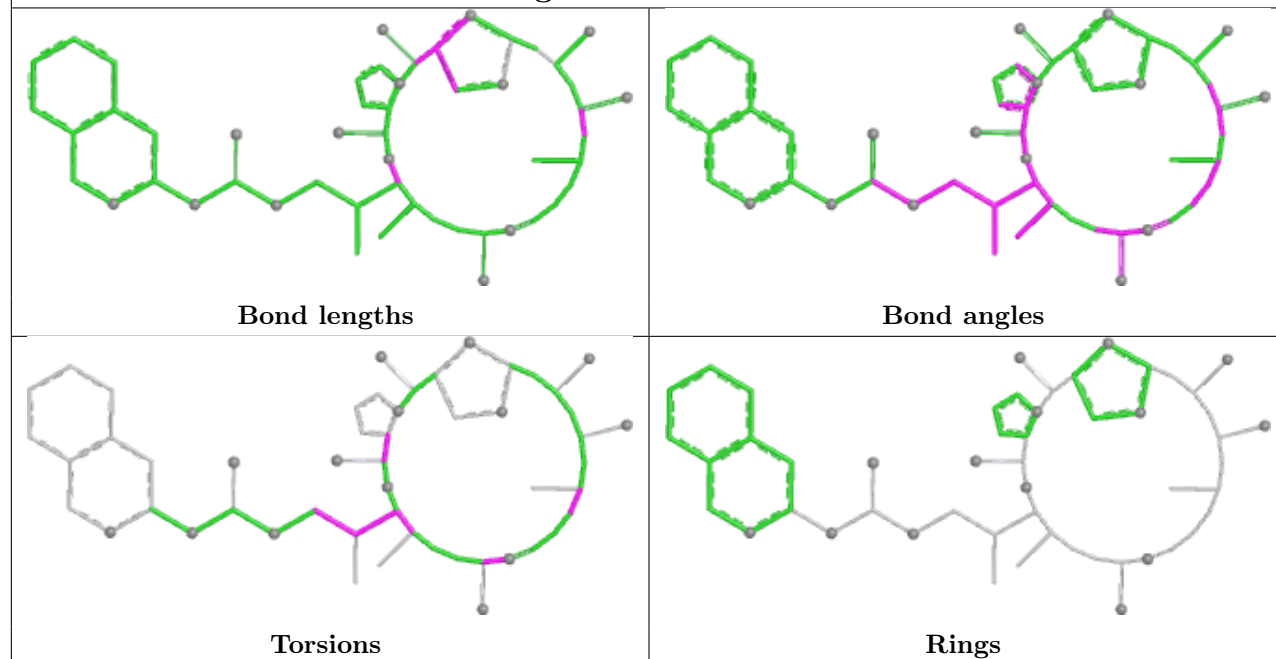
Ligand SXA D 302



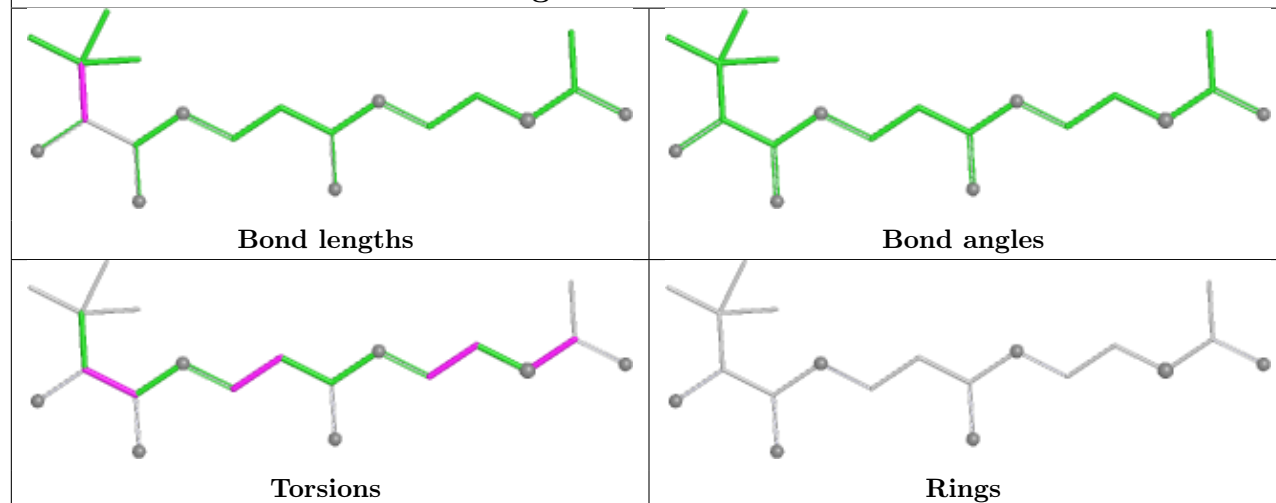
Ligand SXA F 405



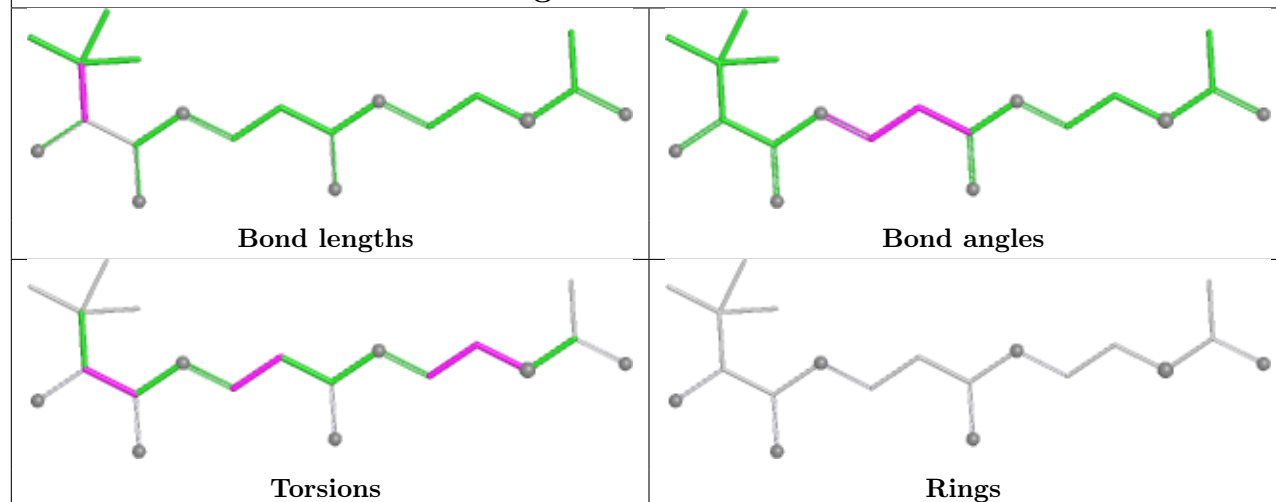
Ligand O7V E 401



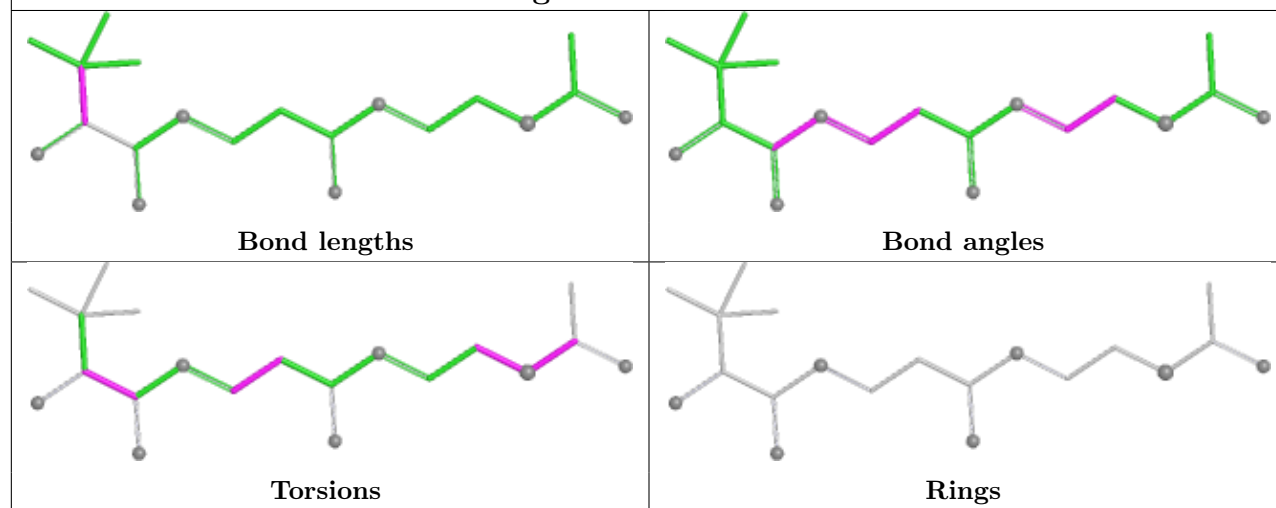
Ligand SXA A 403



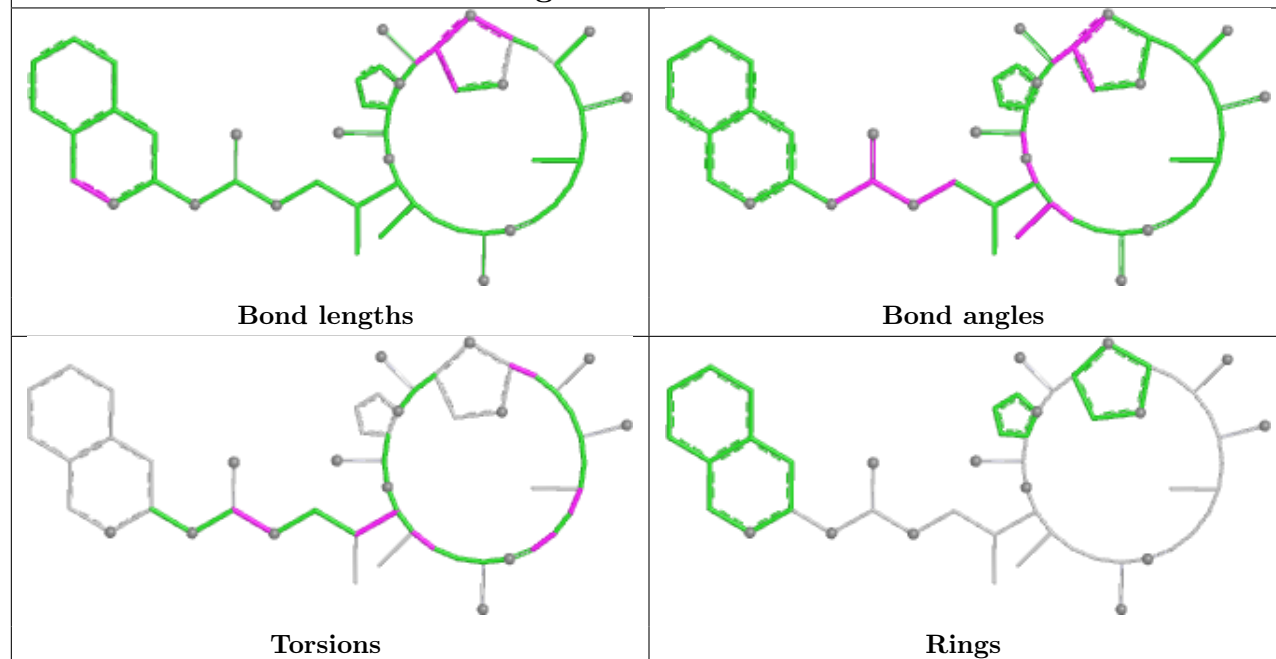
Ligand SXA F 406



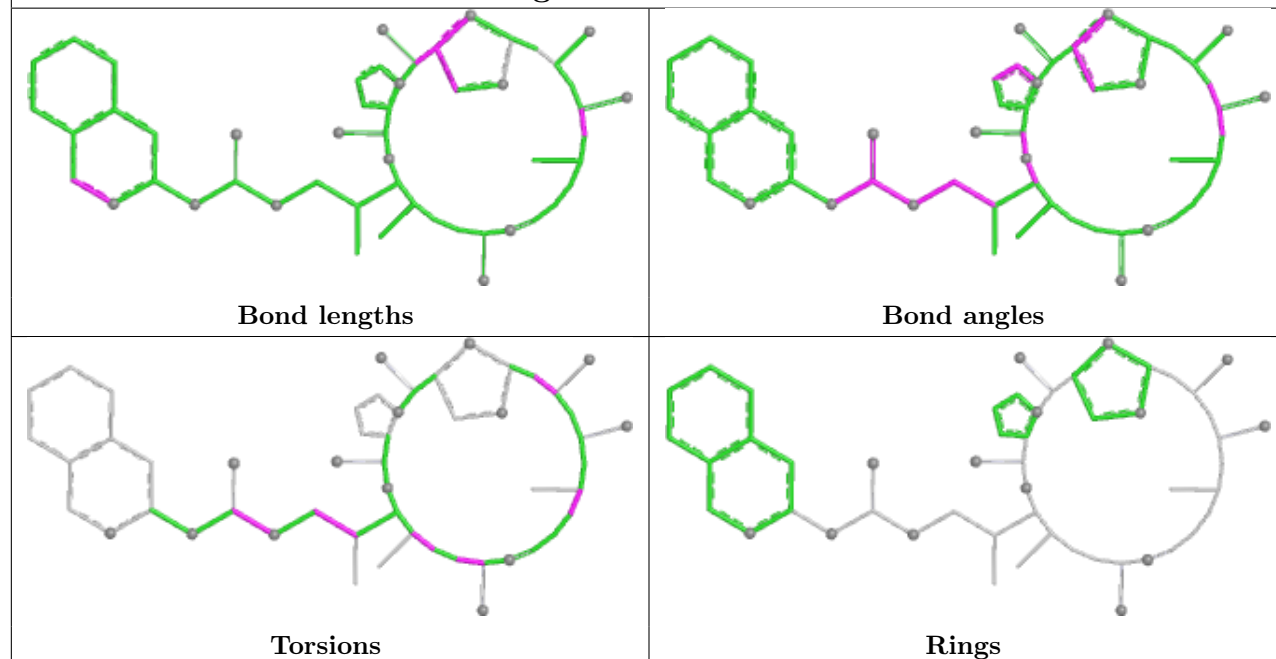
Ligand SXA B 405



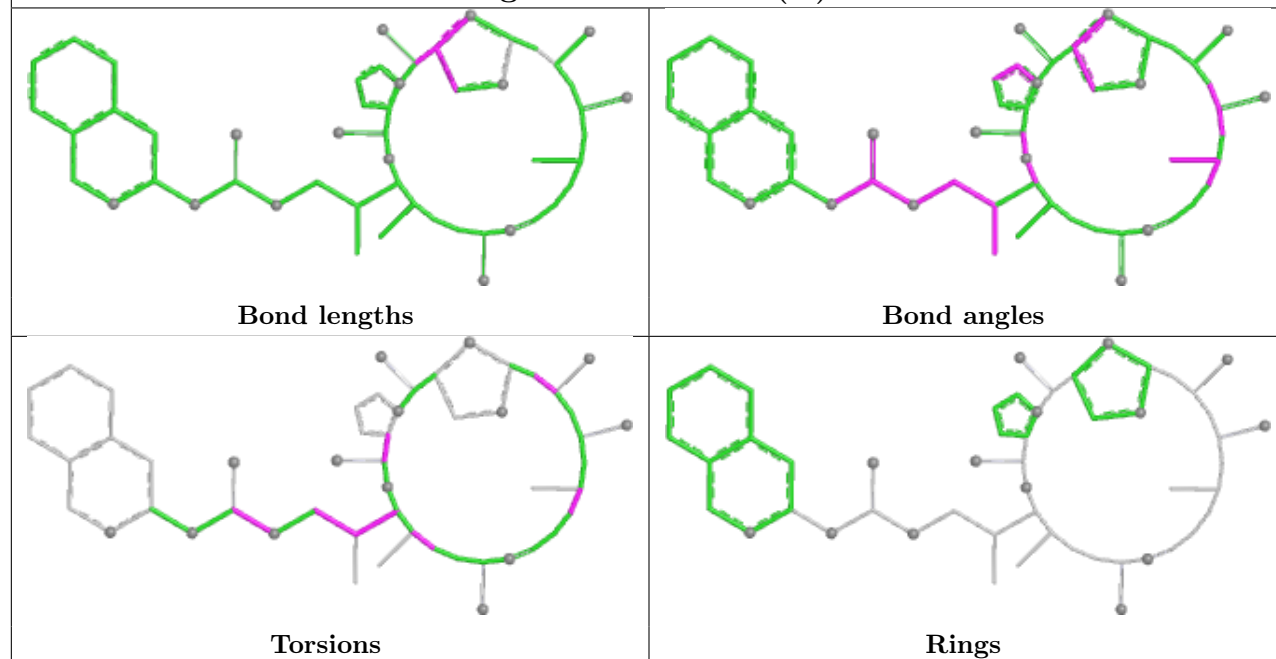
Ligand O7V F 401



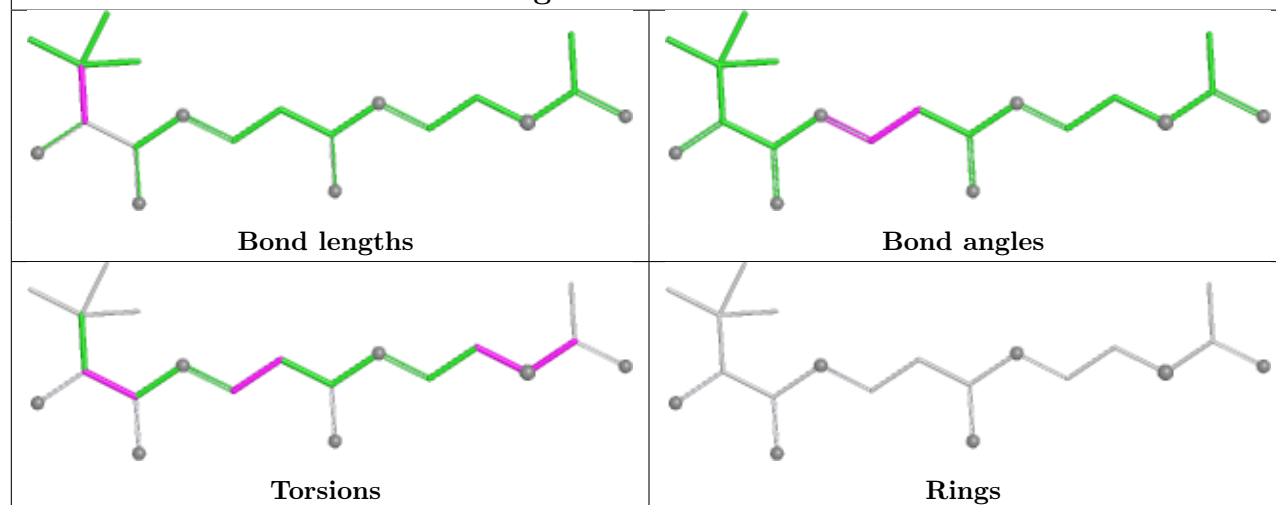
Ligand O7V A 401

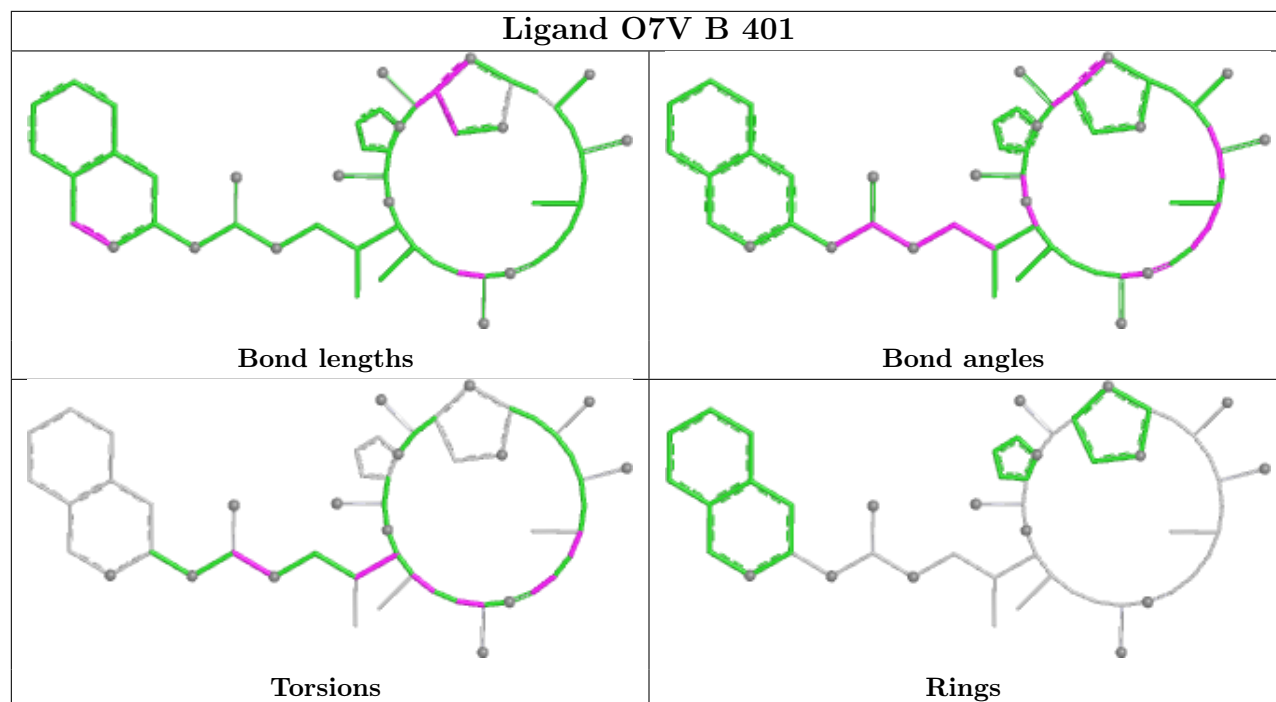


Ligand O7V C 401 (A)



Ligand SXA E 407





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	206/209 (98%)	0.28	10 (4%) 35 31	33, 52, 86, 118	0
1	B	206/209 (98%)	0.45	14 (6%) 23 20	34, 49, 78, 118	0
1	C	209/209 (100%)	0.32	15 (7%) 21 18	30, 51, 93, 113	0
1	D	207/209 (99%)	0.51	17 (8%) 17 15	37, 53, 82, 115	0
1	E	207/209 (99%)	0.54	16 (7%) 19 17	34, 56, 89, 119	0
1	F	204/209 (97%)	0.67	18 (8%) 15 13	32, 54, 91, 121	0
All	All	1239/1254 (98%)	0.46	90 (7%) 21 18	30, 52, 88, 121	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	7	HIS	7.9
1	E	21	ARG	6.5
1	D	7	HIS	6.1
1	F	7	HIS	6.0
1	E	7	HIS	5.8
1	B	32	ASN	5.3
1	F	111	LYS	5.3
1	A	7	HIS	5.3
1	C	215	LEU	5.2
1	B	33	GLU	5.1
1	E	207	GLU	4.6
1	F	21	ARG	4.5
1	F	210	LYS	4.5
1	D	21	ARG	4.5
1	E	213	ARG	4.4
1	C	7	HIS	4.3
1	B	107	MET	4.1
1	D	207	GLU	4.0
1	C	205	ASP	4.0

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Mol	Chain	Res	Type	RSRZ
1	D	210	LYS	3.8
1	D	213	ARG	3.8
1	F	33	GLU	3.6
1	D	211	ARG	3.5
1	A	211	ARG	3.4
1	C	107	MET	3.4
1	A	212	LYS	3.4
1	E	18	LYS	3.4
1	C	91	GLY	3.4
1	D	203	ASN	3.4
1	C	134	MET	3.3
1	C	102	GLY	3.3
1	C	33	GLU	3.3
1	A	203	ASN	3.3
1	F	196	GLU	3.3
1	B	192	LYS	3.3
1	B	21	ARG	3.2
1	B	210	LYS	3.2
1	E	211	ARG	3.2
1	F	209	LEU	3.1
1	F	200	PHE	3.1
1	F	35	ILE	3.0
1	F	39	GLU	2.9
1	E	105	LYS	2.8
1	B	208	MET	2.8
1	D	92	SER	2.8
1	D	209	LEU	2.8
1	A	105	LYS	2.8
1	D	191	MET	2.8
1	E	47	ARG	2.7
1	F	18	LYS	2.7
1	F	192	LYS	2.7
1	C	111	LYS	2.6
1	E	179	GLU	2.6
1	B	193	ILE	2.6
1	E	107	MET	2.6
1	E	192	LYS	2.5
1	F	67	ILE	2.5
1	D	18	LYS	2.5
1	A	47	ARG	2.5
1	A	184	LEU	2.5
1	F	191	MET	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	111	LYS	2.4
1	D	102	GLY	2.4
1	E	106	TYR	2.3
1	D	192	LYS	2.3
1	B	159	SER	2.3
1	C	208	MET	2.2
1	D	107	MET	2.2
1	F	12	GLU	2.2
1	B	47	ARG	2.2
1	C	105	LYS	2.2
1	D	206	ILE	2.2
1	D	212	LYS	2.2
1	F	24	GLN	2.2
1	C	214	LYS	2.2
1	C	167	LYS	2.1
1	E	212	LYS	2.1
1	F	207	GLU	2.1
1	B	31	THR	2.1
1	F	70	ARG	2.1
1	A	111	LYS	2.1
1	A	91	GLY	2.1
1	E	99	PHE	2.1
1	C	47	ARG	2.1
1	A	48	GLY	2.1
1	D	205	ASP	2.0
1	B	207	GLU	2.0
1	B	212	LYS	2.0
1	C	213	ARG	2.0
1	E	100	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

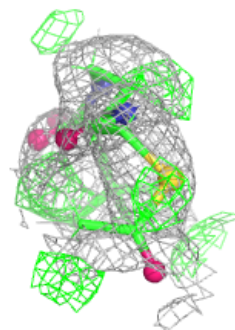
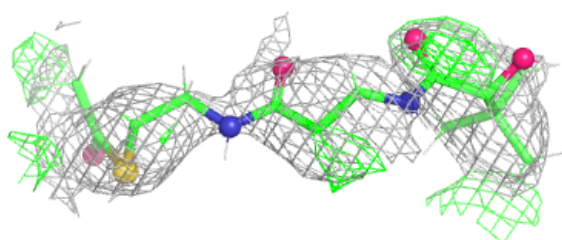
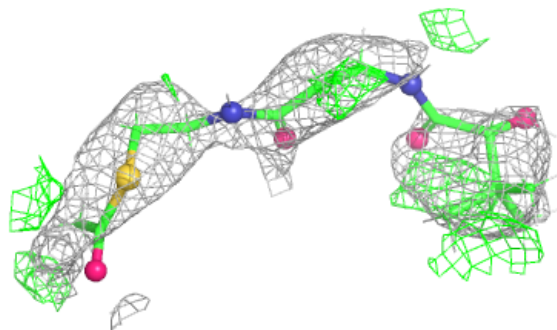
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	PO4	A	402	5/5	0.73	0.26	71,71,71,71	5
4	SXA	F	405	20/25	0.77	0.31	0,0,85,85	7
4	SXA	D	302	20/25	0.78	0.32	0,0,82,82	9
3	PO4	E	406	5/5	0.78	0.26	77,77,78,78	5
4	SXA	B	405	20/25	0.79	0.32	0,0,89,90	7
4	SXA	F	406	20/25	0.79	0.30	0,0,86,86	9
6	MG	D	301	1/1	0.81	0.13	60,60,60,60	0
2	O7V	C	401[B]	52/52	0.82	0.26	0,63,68,69	96
2	O7V	C	401[A]	52/52	0.82	0.26	0,63,68,69	96
4	SXA	A	403	20/25	0.83	0.25	0,0,63,64	11
2	O7V	E	401	52/52	0.83	0.18	0,76,79,79	0
5	CL	F	403	1/1	0.84	0.13	87,87,87,87	0
4	SXA	E	407	20/25	0.85	0.23	0,0,70,71	8
2	O7V	A	401	52/52	0.86	0.17	0,62,66,66	0
2	O7V	F	401	52/52	0.86	0.16	0,54,58,59	0
2	O7V	B	401	52/52	0.88	0.17	0,56,61,61	23
2	O7V	B	402	52/52	0.89	0.16	0,58,60,60	0
5	CL	C	402	1/1	0.91	0.27	66,66,66,66	0
5	CL	E	402	1/1	0.92	0.26	73,73,73,73	0
6	MG	C	404	1/1	0.92	0.09	45,45,45,45	0
5	CL	E	405	1/1	0.92	0.25	91,91,91,91	0
5	CL	E	404	1/1	0.93	0.22	90,90,90,90	0
5	CL	E	403	1/1	0.94	0.09	64,64,64,64	0
5	CL	C	403	1/1	0.94	0.11	85,85,85,85	0
5	CL	F	404	1/1	0.95	0.19	122,122,122,122	0
5	CL	B	403	1/1	0.96	0.09	65,65,65,65	0
5	CL	B	404	1/1	0.97	0.07	57,57,57,57	0
5	CL	F	402	1/1	0.97	0.19	74,74,74,74	0

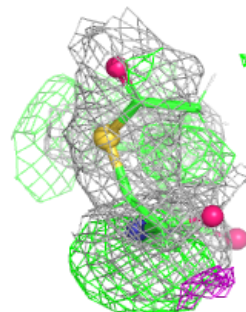
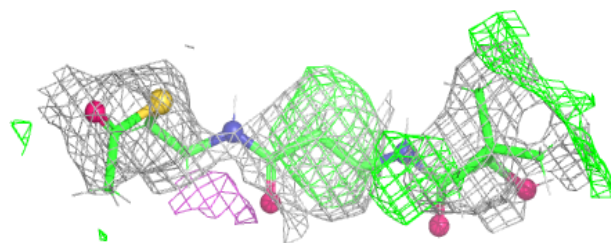
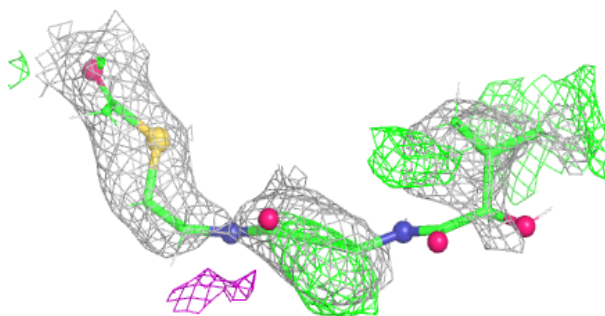
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around SXA F 405:

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

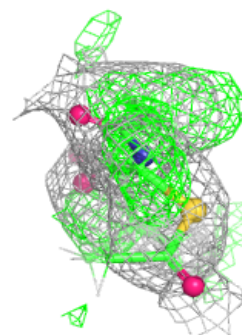
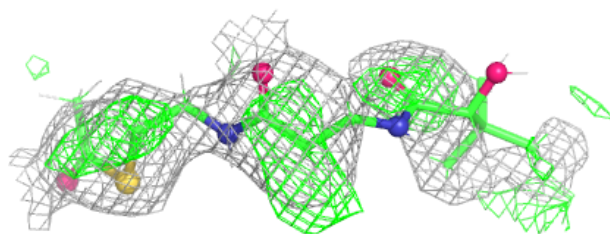
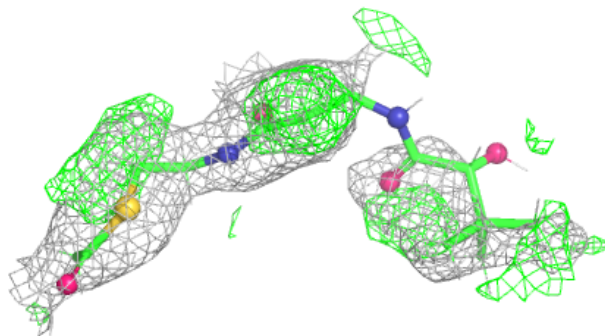
**Electron density around SXA D 302:**

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

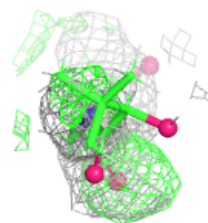
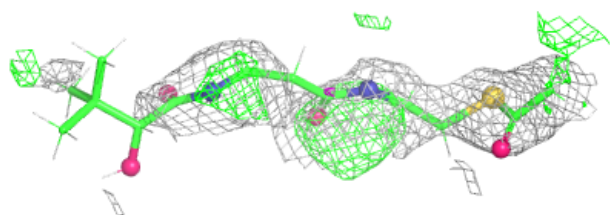
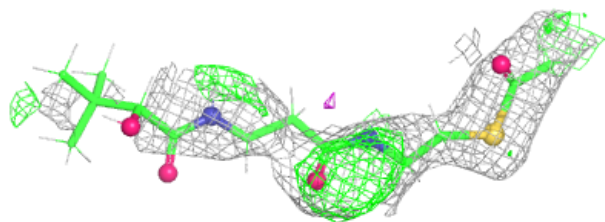


Electron density around SXA B 405:

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

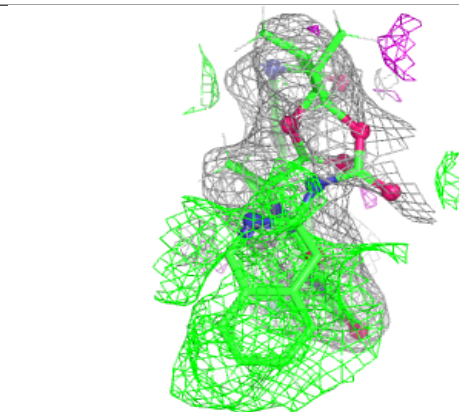
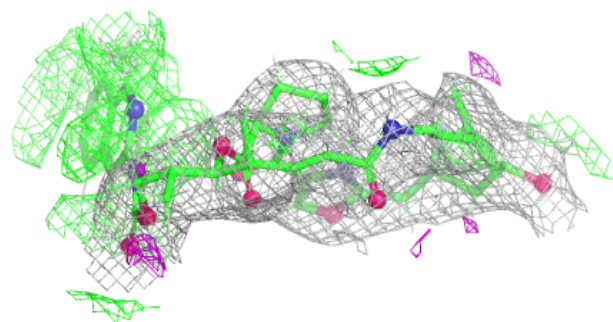
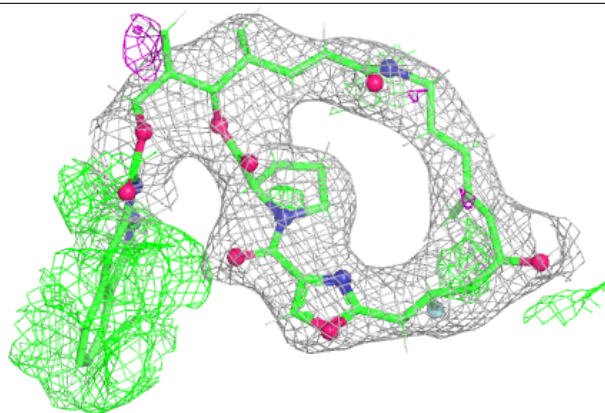
**Electron density around SXA F 406:**

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

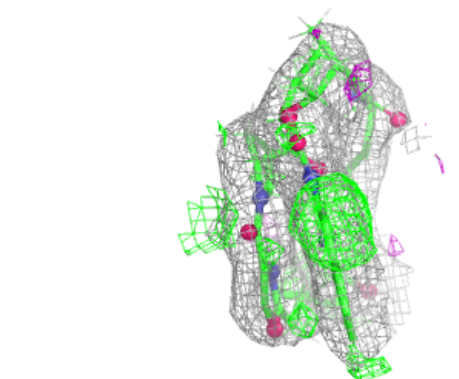
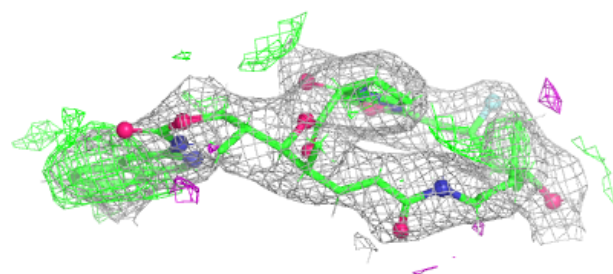
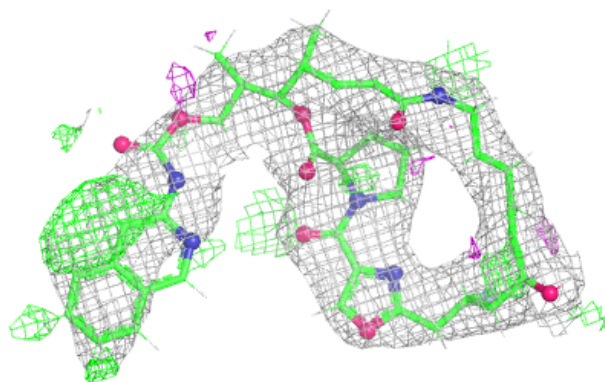


Electron density around O7V C 401 (B):

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

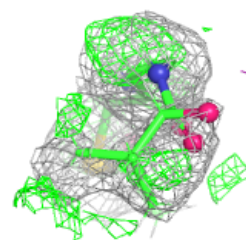
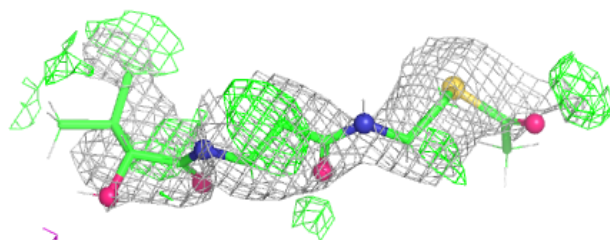
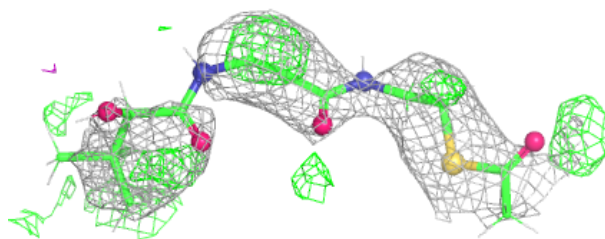
**Electron density around O7V C 401 (A):**

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

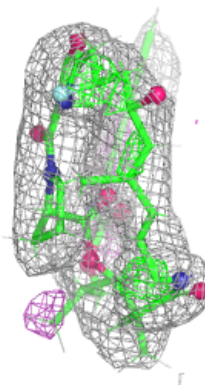
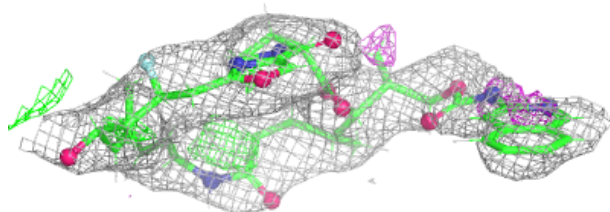
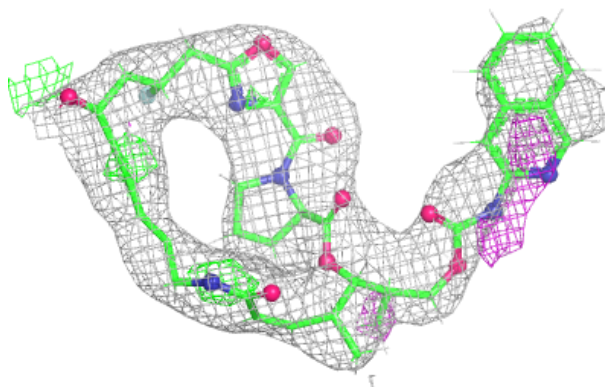


Electron density around SXA A 403:

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

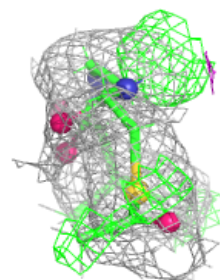
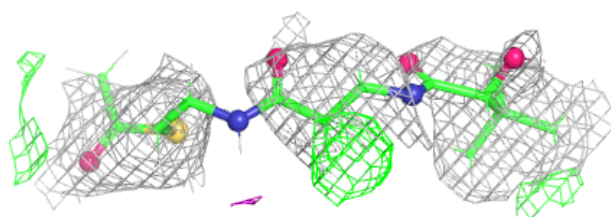
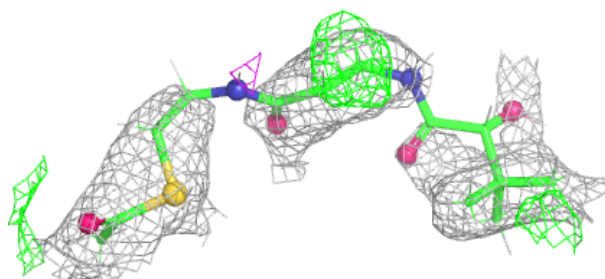
**Electron density around O7V E 401:**

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

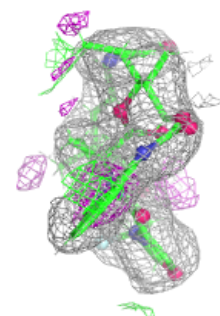
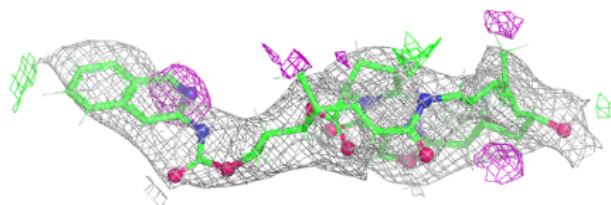
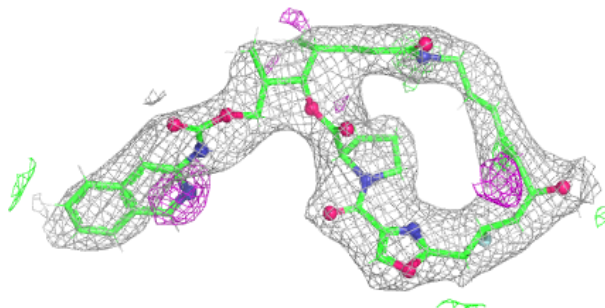


Electron density around SXA E 407:

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

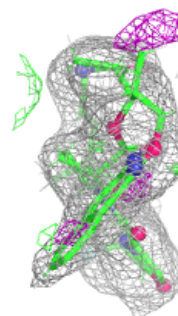
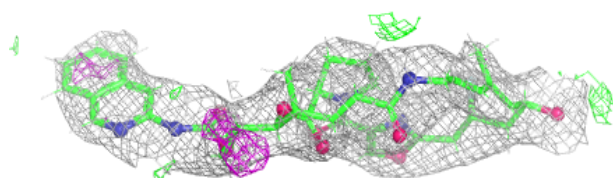
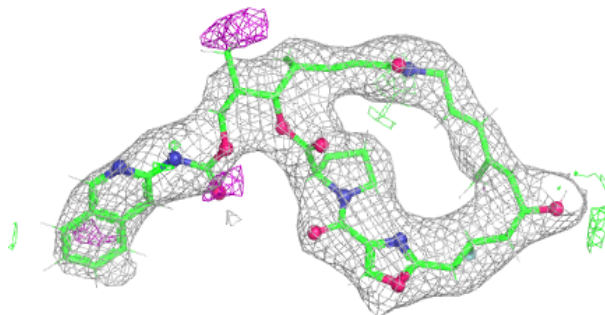
**Electron density around O7V A 401:**

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

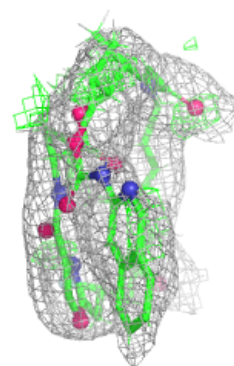
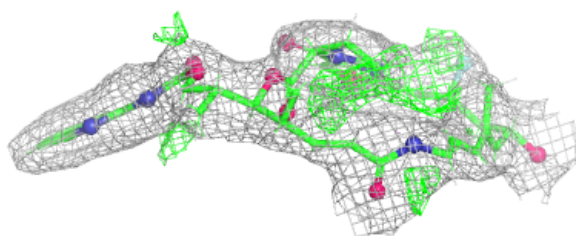
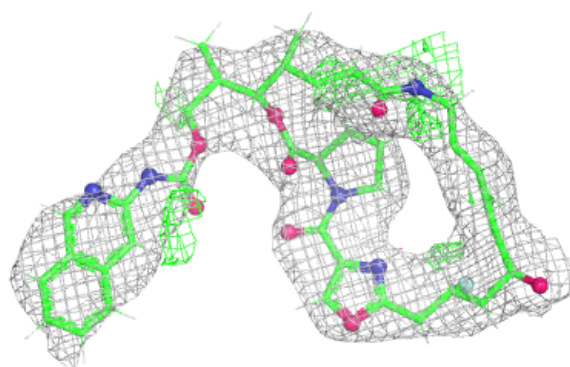


Electron density around O7V F 401:

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

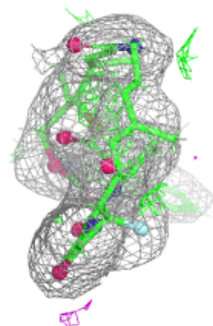
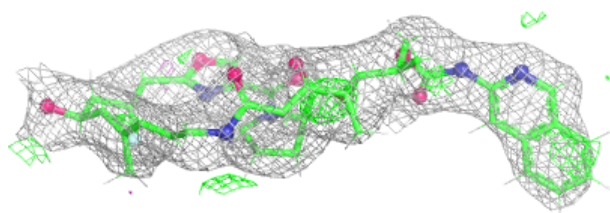
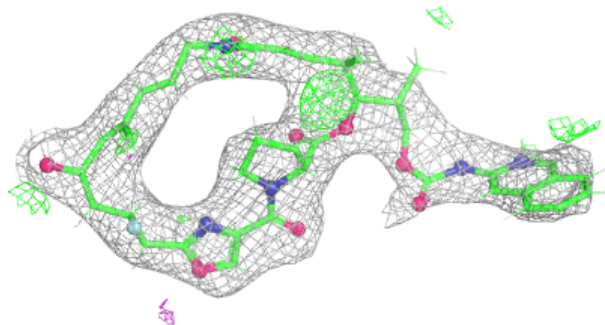
**Electron density around O7V B 401:**

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



Electron density around O7V B 402:

$2mF_o-DF_c$ (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.