



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:07 PM UTC

PDB ID : 4KF6 / pdb_00004kf6
Title : Crystal structure of human ceramide-1-phosphate transfer protein (CPTP) in complex with 8:0 Ceramide-1-Phosphate (8:0-C1P)
Authors : Simanshu, D.K.; Brown, R.E.; Patel, D.J.
Deposited on : 2013-04-26
Resolution : 3.19 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

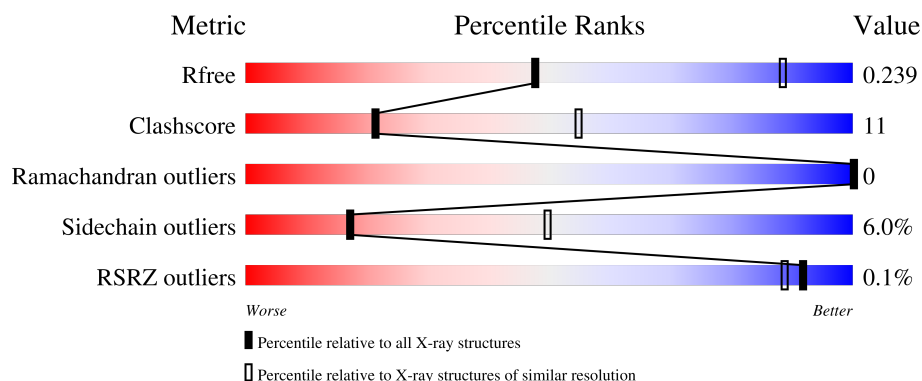
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.19 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	215	
1	B	215	
1	C	215	
1	D	215	
1	E	215	

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Mol	Chain	Length	Quality of chain
1	F	215	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: green (61%), yellow (30%), and orange (5%). The segments are labeled with their respective percentages: 61%, 30%, and 5%. The bar is set against a background of a light gray grid.

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9915 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.

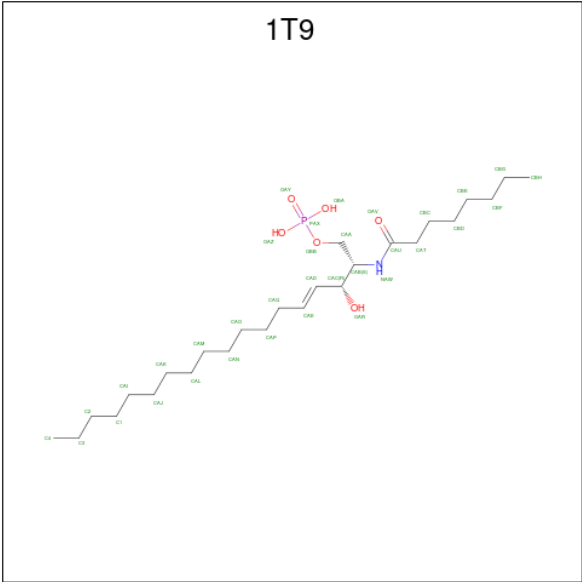
- Molecule 1 is a protein called Glycolipid transfer protein domain-containing protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	206	Total	C	N	O	S	0	0	0
			1625	1035	291	291	8			
1	B	208	Total	C	N	O	S	0	0	0
			1647	1050	296	293	8			
1	C	206	Total	C	N	O	S	0	0	0
			1640	1047	293	292	8			
1	D	206	Total	C	N	O	S	0	0	0
			1630	1042	291	289	8			
1	E	206	Total	C	N	O	S	0	0	0
			1617	1029	293	287	8			
1	F	206	Total	C	N	O	S	0	0	0
			1598	1016	287	287	8			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q5TA50
B	0	SER	-	expression tag	UNP Q5TA50
C	0	SER	-	expression tag	UNP Q5TA50
D	0	SER	-	expression tag	UNP Q5TA50
E	0	SER	-	expression tag	UNP Q5TA50
F	0	SER	-	expression tag	UNP Q5TA50

- Molecule 2 is (2S,3R,4E)-3-hydroxy-2-(octanoylamino)octadec-4-en-1-yl dihydrogen phosphate (CCD ID: 1T9) (formula: C₂₆H₅₂NO₆P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			30	22	1	6	1		
2	B	1	Total	C	N	O	P	0	0
			25	17	1	6	1		
2	C	1	Total	C	N	O	P	0	0
			29	21	1	6	1		
2	D	1	Total	C	N	O	P	0	0
			27	19	1	6	1		
2	E	1	Total	C	N	O	P	0	0
			22	14	1	6	1		
2	F	1	Total	C	N	O	P	0	0
			16	8	1	6	1		

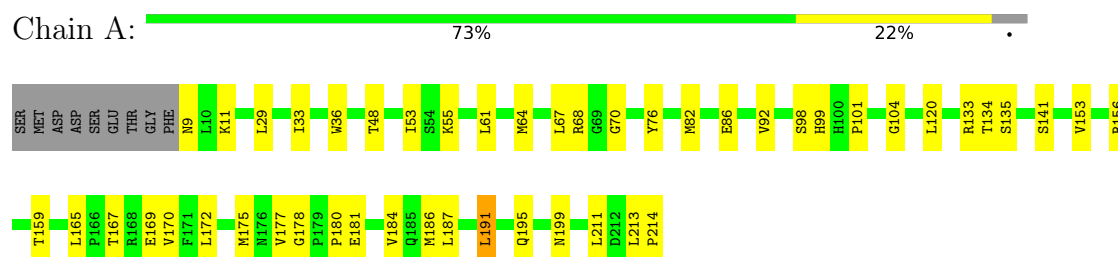
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total	O	0	0
			4	4		
3	B	2	Total	O	0	0
			2	2		
3	D	1	Total	O	0	0
			1	1		
3	E	2	Total	O	0	0
			2	2		

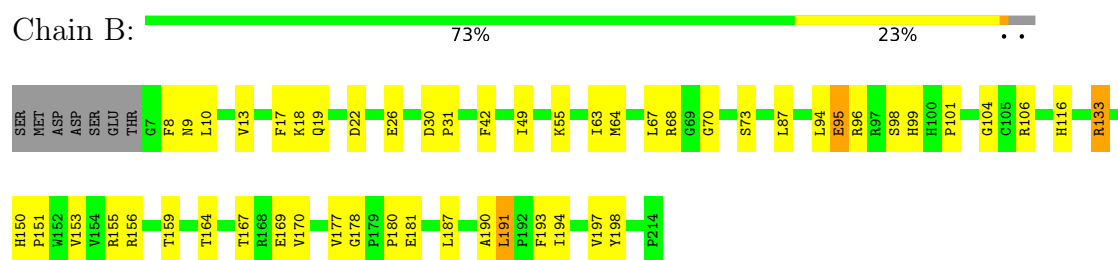
3 Residue-property plots [i](#)

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.

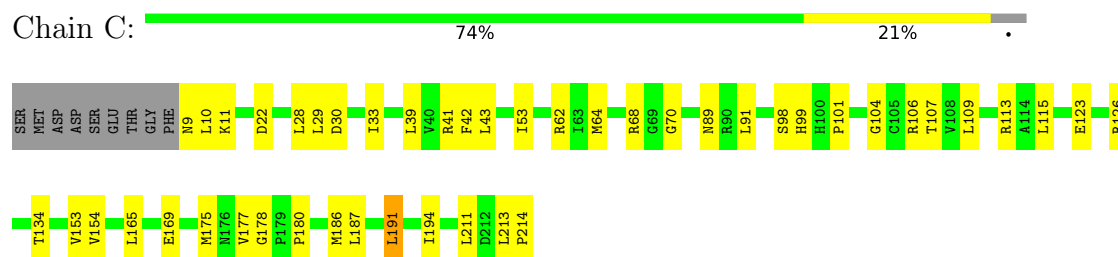
- Molecule 1: Glycolipid transfer protein domain-containing protein 1



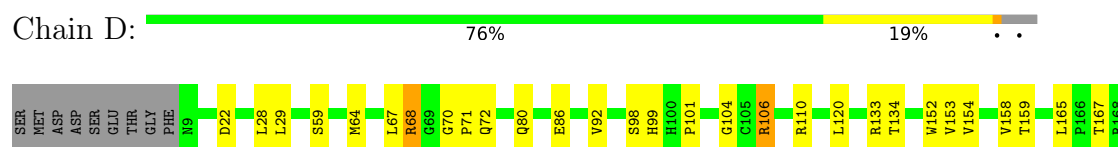
- Molecule 1: Glycolipid transfer protein domain-containing protein 1



- Molecule 1: Glycolipid transfer protein domain-containing protein 1

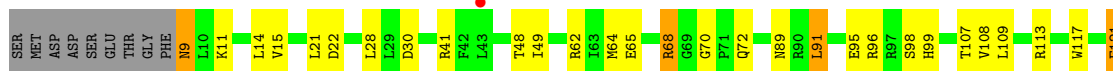


- Molecule 1: Glycolipid transfer protein domain-containing protein 1





- Molecule 1: Glycolipid transfer protein domain-containing protein 1



- Molecule 1: Glycolipid transfer protein domain-containing protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.88Å 122.40Å 149.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.83 – 3.19 38.83 – 3.19	Depositor EDS
% Data completeness (in resolution range)	99.8 (38.83-3.19) 99.7 (38.83-3.19)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 3.18Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.222 , 0.265 (Not available) , 0.239	Depositor DCC
R_{free} test set	1302 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	83.3	Xtriage
Anisotropy	0.138	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 68.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9915	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.41% 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: 1T9

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	0.42	0/1661	0.90	5/2255 (0.2%)
1	B	0.44	0/1685	0.91	5/2285 (0.2%)
1	C	0.37	0/1677	0.84	4/2275 (0.2%)
1	D	0.39	0/1667	0.87	4/2263 (0.2%)
1	E	0.40	0/1654	0.86	4/2246 (0.2%)
1	F	0.49	1/1634 (0.1%)	1.01	5/2220 (0.2%)
All	All	0.42	1/9978 (0.0%)	0.90	27/13544 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	85	HIS	N-CA	5.56	1.53	1.46

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	177	VAL	N-CA-C	-7.40	106.10	113.20
1	B	177	VAL	N-CA-C	-6.83	106.65	113.20
1	A	178	GLY	CA-C-N	6.75	127.34	120.38
1	A	178	GLY	C-N-CA	6.75	127.34	120.38
1	D	178	GLY	CA-C-N	6.73	127.31	120.38
1	D	178	GLY	C-N-CA	6.73	127.31	120.38
1	A	70	GLY	CA-C-N	6.67	127.20	119.47
1	A	70	GLY	C-N-CA	6.67	127.20	119.47
1	B	178	GLY	CA-C-N	6.18	126.75	120.38
1	B	178	GLY	C-N-CA	6.18	126.75	120.38
1	E	178	GLY	CA-C-N	6.18	126.75	120.38
1	E	178	GLY	C-N-CA	6.18	126.75	120.38
1	C	178	GLY	CA-C-N	5.98	126.53	120.38
1	C	178	GLY	C-N-CA	5.98	126.53	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	70	GLY	CA-C-N	5.95	126.37	119.47
1	D	70	GLY	C-N-CA	5.95	126.37	119.47
1	B	70	GLY	CA-C-N	5.92	127.24	119.84
1	B	70	GLY	C-N-CA	5.92	127.24	119.84
1	F	85	HIS	CA-C-O	-5.84	114.68	120.70
1	F	70	GLY	CA-C-N	5.80	127.09	119.84
1	F	70	GLY	C-N-CA	5.80	127.09	119.84
1	E	70	GLY	CA-C-N	5.67	126.05	119.47
1	E	70	GLY	C-N-CA	5.67	126.05	119.47
1	C	70	GLY	CA-C-N	5.66	126.03	119.47
1	C	70	GLY	C-N-CA	5.66	126.03	119.47
1	F	128	SER	CA-C-N	5.44	125.74	119.92
1	F	128	SER	C-N-CA	5.44	125.74	119.92

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	1625	0	1626	29	0
1	B	1647	0	1640	33	0
1	C	1640	0	1651	33	0
1	D	1630	0	1638	22	0
1	E	1617	0	1600	43	0
1	F	1598	0	1569	59	0
2	A	30	0	39	4	0
2	B	25	0	26	0	0
2	C	29	0	37	6	0
2	D	27	0	30	2	0
2	E	22	0	23	1	0
2	F	16	0	8	2	0
3	A	4	0	0	0	0
3	B	2	0	0	1	0
3	D	1	0	0	0	0
3	E	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	9915	0	9887	215	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 11.

All (215) 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:18:LYS:HB2	1:B:193:PHE:HZ	1.29	0.96
1:F:75:HIS:NE2	1:F:85:HIS:HB3	1.86	0.88
1:F:75:HIS:CD2	1:F:85:HIS:HB3	2.09	0.87
1:C:53:ILE:HG21	2:C:301:1T9:H40	1.55	0.86
1:D:29:LEU:HD22	1:D:68:ARG:HG3	1.57	0.85
1:F:132:ALA:N	1:F:168:ARG:HH22	1.78	0.82
1:F:64:MET:HE1	1:F:104:GLY:HA2	1.63	0.81
1:D:64:MET:HE1	1:D:104:GLY:HA2	1.61	0.80
1:E:64:MET:HE3	1:E:108:VAL:HG23	1.62	0.80
1:C:64:MET:HE1	1:C:104:GLY:HA2	1.64	0.79
1:A:64:MET:HE1	1:A:104:GLY:HA2	1.64	0.77
1:C:53:ILE:CG2	2:C:301:1T9:H40	2.15	0.77
1:E:30:ASP:OD2	1:E:68:ARG:NH1	2.20	0.75
1:F:21:LEU:HD21	1:F:200:VAL:HG11	1.69	0.74
1:E:64:MET:HE2	1:E:107:THR:HB	1.69	0.73
1:E:21:LEU:HD21	1:E:200:VAL:HG11	1.69	0.73
1:B:64:MET:HE1	1:B:104:GLY:HA2	1.70	0.72
1:F:129:PRO:O	1:F:168:ARG:NH2	2.25	0.70
1:E:195:GLN:O	1:E:199:ASN:ND2	2.24	0.69
1:B:95:GLU:N	1:B:95:GLU:OE1	2.24	0.68
1:D:110:ARG:HG2	2:D:301:1T9:H26	1.75	0.67
1:C:169:GLU:HG3	1:C:180:PRO:HG3	1.77	0.67
1:F:167:THR:HG22	1:F:170:VAL:HG12	1.77	0.66
1:F:177:VAL:HG21	1:F:186:MET:SD	2.37	0.65
1:F:134:THR:HG23	1:F:165:LEU:HD13	1.79	0.64
1:F:135:SER:HA	1:F:165:LEU:HD21	1.78	0.64
1:B:68:ARG:NH1	1:B:73:SER:O	2.31	0.64
1:C:9:ASN:HD21	1:C:41:ARG:HH12	1.46	0.63
1:F:168:ARG:HG3	1:F:169:GLU:H	1.64	0.62
1:B:133:ARG:NH2	3:B:402:HOH:O	2.32	0.62
1:F:132:ALA:H	1:F:168:ARG:HH22	1.48	0.61
1:E:192:PRO:O	1:E:196:ARG:HG2	1.99	0.61
1:B:18:LYS:HB2	1:B:193:PHE:CZ	2.22	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:95:GLU:HG3	1:E:96:ARG:H	1.67	0.59
1:D:169:GLU:HG3	1:D:180:PRO:HG3	1.85	0.59
1:C:30:ASP:OD2	1:C:68:ARG:NH1	2.35	0.59
1:D:133:ARG:NH2	1:D:165:LEU:O	2.29	0.59
1:E:9:ASN:HD22	1:E:41:ARG:HH12	1.50	0.58
1:E:182:GLN:OE1	1:E:185:GLN:NE2	2.36	0.58
1:D:179:PRO:HD2	1:D:182:GLN:OE1	2.03	0.58
1:E:167:THR:OG1	1:E:169:GLU:OE2	2.21	0.58
1:A:67:LEU:HD23	1:A:76:TYR:CZ	2.39	0.57
1:E:152:TRP:HE1	2:F:301:1T9:H21	1.68	0.57
1:A:36:TRP:HB3	1:A:61:LEU:HD11	1.87	0.57
1:F:19:GLN:OE1	1:F:19:GLN:N	2.38	0.56
1:A:172:LEU:HA	1:A:175:MET:HE3	1.87	0.56
1:F:36:TRP:CE3	1:F:39:LEU:HD11	2.41	0.56
1:F:106:ARG:NH1	1:F:214:PRO:OXT	2.39	0.56
1:A:134:THR:HG22	1:A:165:LEU:HD13	1.88	0.55
1:C:115:LEU:HD21	1:C:194:ILE:HG12	1.88	0.55
1:E:62:ARG:NH1	1:E:65:GLU:OE1	2.39	0.55
1:D:167:THR:HG22	1:D:170:VAL:HG23	1.89	0.55
1:F:115:LEU:HD11	1:F:194:ILE:HG21	1.89	0.55
1:A:98:SER:OG	1:A:99:HIS:N	2.36	0.55
1:E:22:ASP:HB3	1:E:28:LEU:HD11	1.88	0.54
1:B:87:LEU:HD21	1:B:94:LEU:HD21	1.89	0.54
1:C:98:SER:OG	1:C:99:HIS:N	2.28	0.54
1:A:55:LYS:HD2	1:E:48:THR:HG21	1.89	0.54
1:D:110:ARG:CG	2:D:301:1T9:H26	2.38	0.53
1:A:11:LYS:HE2	1:A:186:MET:HG3	1.90	0.53
1:E:134:THR:HG22	1:E:165:LEU:HD13	1.90	0.53
1:A:195:GLN:O	1:A:199:ASN:ND2	2.40	0.53
1:E:185:GLN:HG3	1:E:186:MET:N	2.23	0.53
1:D:98:SER:OG	1:D:99:HIS:ND1	2.42	0.53
1:F:75:HIS:HE1	1:F:86:GLU:OE1	1.92	0.53
1:F:86:GLU:HB3	1:F:92:VAL:HG23	1.89	0.53
1:F:44:ASN:H	1:F:44:ASN:HD22	1.56	0.53
1:A:48:THR:OG1	1:B:156:ARG:NH2	2.42	0.52
1:B:181:GLU:OE1	1:B:181:GLU:N	2.36	0.52
1:C:22:ASP:HB3	1:C:28:LEU:HD11	1.91	0.52
1:F:192:PRO:O	1:F:196:ARG:HG3	2.09	0.52
1:A:33:ILE:HG23	1:A:61:LEU:HD22	1.92	0.52
1:B:30:ASP:CG	1:B:68:ARG:HE	2.17	0.52
1:B:98:SER:OG	1:B:99:HIS:N	2.41	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:ARG:NH1	1:A:135:SER:H	2.08	0.52
1:B:190:ALA:O	1:B:193:PHE:HB3	2.10	0.52
1:C:187:LEU:O	1:C:191:LEU:HB2	2.10	0.52
1:D:133:ARG:HH11	1:D:134:THR:H	1.57	0.52
1:F:17:PHE:CE2	1:F:35:SER:HB3	2.45	0.52
1:F:115:LEU:HD11	1:F:194:ILE:CG2	2.39	0.52
1:A:9:ASN:ND2	1:A:9:ASN:N	2.58	0.52
1:E:121:PHE:CD2	1:E:121:PHE:C	2.87	0.52
1:F:183:ALA:HA	1:F:186:MET:HB3	1.91	0.51
1:F:194:ILE:N	1:F:194:ILE:HD13	2.24	0.51
1:C:106:ARG:NH1	1:C:214:PRO:OXT	2.42	0.51
1:F:162:PHE:O	1:F:165:LEU:HG	2.10	0.51
1:D:99:HIS:O	1:D:101:PRO:HD3	2.10	0.51
1:B:167:THR:HG22	1:B:170:VAL:HG22	1.92	0.51
1:D:98:SER:OG	1:D:99:HIS:N	2.43	0.51
1:F:98:SER:OG	1:F:99:HIS:N	2.42	0.51
1:A:76:TYR:CE1	1:A:82:MET:HG3	2.46	0.50
1:C:134:THR:HG22	1:C:165:LEU:HD13	1.92	0.50
1:E:121:PHE:HE1	1:E:138:CYS:SG	2.33	0.50
1:F:135:SER:HB2	1:F:165:LEU:HD11	1.94	0.50
1:C:9:ASN:HD21	1:C:41:ARG:NH1	2.09	0.50
1:F:36:TRP:O	1:F:40:VAL:HG23	2.12	0.50
1:F:60:LYS:NZ	2:F:301:1T9:H26	2.25	0.50
1:E:191:LEU:N	1:E:192:PRO:HD2	2.26	0.50
1:F:169:GLU:HG3	1:F:180:PRO:HG3	1.93	0.49
1:C:123:GLU:CD	1:C:126:ARG:HH21	2.20	0.49
1:B:19:GLN:O	1:B:31:PRO:HG2	2.12	0.49
1:A:181:GLU:OE1	1:A:181:GLU:N	2.33	0.49
1:E:15:VAL:HA	1:E:193:PHE:CD2	2.47	0.49
1:A:156:ARG:NH2	2:C:301:1T9:OAY	2.41	0.49
1:F:75:HIS:CE1	1:F:82:MET:HA	2.48	0.49
1:A:187:LEU:O	1:A:191:LEU:HB2	2.13	0.49
1:B:49:ILE:HG21	1:B:164:THR:HG21	1.95	0.49
1:D:22:ASP:HB3	1:D:28:LEU:HD11	1.94	0.48
1:F:56:ASP:O	1:F:59:SER:OG	2.26	0.48
1:B:150:HIS:O	1:B:155:ARG:NH1	2.46	0.48
1:F:79:LEU:O	1:F:82:MET:HB3	2.14	0.48
1:F:17:PHE:CD2	1:F:194:ILE:HD12	2.49	0.48
1:F:167:THR:HG23	1:F:168:ARG:HG3	1.95	0.48
1:F:75:HIS:CD2	1:F:85:HIS:CB	2.92	0.48
1:B:68:ARG:HG3	1:B:68:ARG:HH11	1.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:GLU:CD	1:B:96:ARG:H	2.22	0.47
1:A:53:ILE:HG23	2:A:301:1T9:H38	1.96	0.47
1:D:106:ARG:NH1	1:D:214:PRO:OXT	2.43	0.47
1:D:133:ARG:NH2	1:E:170:VAL:HG21	2.29	0.47
1:A:120:LEU:HB3	1:A:141:SER:HB2	1.96	0.47
1:E:89:ASN:HB2	1:E:91:LEU:HD22	1.96	0.47
1:E:11:LYS:HZ3	1:E:187:LEU:HD23	1.80	0.47
1:F:99:HIS:O	1:F:101:PRO:HD3	2.15	0.47
1:F:108:VAL:O	1:F:112:HIS:N	2.35	0.47
1:E:169:GLU:HG2	1:E:170:VAL:N	2.30	0.46
1:E:49:ILE:HD13	1:E:161:ALA:HA	1.97	0.46
1:F:36:TRP:HB3	1:F:61:LEU:HD21	1.96	0.46
1:C:106:ARG:NH2	2:C:301:1T9:OAZ	2.49	0.46
1:C:29:LEU:O	1:C:33:ILE:HG13	2.16	0.46
1:E:30:ASP:OD1	1:E:68:ARG:HD3	2.15	0.46
1:C:10:LEU:HD13	1:C:42:PHE:CD1	2.51	0.46
1:E:15:VAL:HA	1:E:193:PHE:HD2	1.80	0.46
1:E:98:SER:OG	1:E:99:HIS:N	2.37	0.46
1:A:169:GLU:HG2	1:A:170:VAL:N	2.31	0.46
1:F:119:GLN:HE21	1:F:195:GLN:HE21	1.64	0.45
1:A:153:VAL:HG22	2:A:301:1T9:H4	1.98	0.45
1:B:17:PHE:HB3	1:B:197:VAL:HG21	1.99	0.45
1:B:116:HIS:HB2	1:B:198:TYR:CE1	2.52	0.44
1:F:74:GLU:H	1:F:74:GLU:CD	2.25	0.44
1:A:180:PRO:O	1:A:184:VAL:HG23	2.17	0.44
1:F:134:THR:O	1:F:138:CYS:N	2.44	0.44
1:A:86:GLU:HB3	1:A:92:VAL:HG23	1.99	0.44
1:E:14:LEU:HD21	1:E:122:LEU:HD13	2.00	0.44
1:A:29:LEU:HD23	1:A:68:ARG:HG3	2.00	0.44
1:A:99:HIS:O	1:A:101:PRO:HD3	2.17	0.44
2:A:301:1T9:H9	1:D:152:TRP:CH2	2.53	0.44
1:B:99:HIS:O	1:B:101:PRO:HD3	2.18	0.44
1:F:142:TYR:OH	1:F:150:HIS:ND1	2.32	0.43
1:B:8:PHE:CE2	1:B:42:PHE:HB2	2.53	0.43
1:D:86:GLU:HB3	1:D:92:VAL:HG23	2.00	0.43
1:E:11:LYS:NZ	1:E:187:LEU:HD23	2.32	0.43
1:A:53:ILE:CG2	2:A:301:1T9:H38	2.48	0.43
1:C:62:ARG:HD3	1:C:62:ARG:HA	1.69	0.43
1:C:123:GLU:OE2	1:C:126:ARG:NH2	2.50	0.43
1:B:169:GLU:HG3	1:B:180:PRO:CG	2.49	0.43
1:C:213:LEU:HA	1:C:214:PRO:HD3	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:TYR:HE1	1:A:82:MET:HG3	1.84	0.43
1:B:55:LYS:HE3	1:B:55:LYS:HB2	1.92	0.43
1:E:11:LYS:HD2	1:E:186:MET:HB3	2.01	0.43
1:E:191:LEU:HD12	1:E:191:LEU:HA	1.90	0.43
1:B:22:ASP:OD2	1:B:26:GLU:HG2	2.19	0.43
1:B:150:HIS:HA	1:B:151:PRO:HD3	1.87	0.43
1:C:11:LYS:HE2	1:C:186:MET:HG2	2.00	0.43
1:C:39:LEU:O	1:C:43:LEU:HG	2.19	0.43
1:D:71:PRO:HB2	1:D:72:GLN:OE1	2.18	0.43
1:E:121:PHE:C	1:E:121:PHE:HD2	2.27	0.43
1:E:154:VAL:O	1:E:158:VAL:HG23	2.18	0.43
1:B:63:ILE:O	1:B:67:LEU:HG	2.19	0.43
1:D:154:VAL:O	1:D:158:VAL:HG23	2.19	0.42
1:F:120:LEU:HB3	1:F:141:SER:HB2	2.01	0.42
1:D:80:GLN:HG3	1:D:204:LEU:HD13	2.02	0.42
1:D:174:ALA:O	1:D:176:ASN:ND2	2.52	0.42
1:F:99:HIS:C	1:F:101:PRO:HD3	2.44	0.42
1:F:110:ARG:HD2	1:F:110:ARG:HA	1.88	0.42
1:B:10:LEU:HD12	1:B:13:VAL:CG2	2.49	0.42
1:B:191:LEU:O	1:B:194:ILE:HG22	2.19	0.42
1:A:9:ASN:N	1:A:9:ASN:HD22	2.17	0.42
1:C:106:ARG:HG3	1:C:107:THR:N	2.34	0.42
1:F:36:TRP:CZ3	1:F:39:LEU:HD11	2.54	0.42
1:F:42:PHE:O	1:F:45:SER:OG	2.27	0.42
1:E:109:LEU:O	1:E:113:ARG:HG3	2.19	0.42
1:E:117:TRP:CZ3	2:E:301:1T9:H39	2.55	0.42
1:B:187:LEU:O	1:B:191:LEU:HB2	2.20	0.42
1:E:89:ASN:O	1:E:91:LEU:HD13	2.20	0.42
1:B:169:GLU:HG3	1:B:180:PRO:HG2	2.01	0.41
1:C:153:VAL:CG2	2:C:301:1T9:CAJ	2.98	0.41
1:E:64:MET:HE2	1:E:107:THR:CB	2.45	0.41
1:C:109:LEU:O	1:C:113:ARG:HG3	2.20	0.41
1:E:9:ASN:ND2	1:E:41:ARG:HH12	2.16	0.41
1:C:187:LEU:HD23	1:C:187:LEU:HA	1.94	0.41
1:C:191:LEU:O	1:C:194:ILE:HG22	2.20	0.41
1:F:40:VAL:O	1:F:44:ASN:ND2	2.48	0.41
1:F:72:GLN:HE21	1:F:72:GLN:HB3	1.70	0.41
1:C:154:VAL:HG13	2:C:301:1T9:H16	2.02	0.41
1:F:39:LEU:HD12	1:F:40:VAL:N	2.36	0.41
1:C:169:GLU:HG3	1:C:180:PRO:CG	2.49	0.41
1:E:213:LEU:HA	1:E:214:PRO:HD3	1.82	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:17:PHE:HA	1:F:20:CYS:SG	2.60	0.41
1:B:9:ASN:O	1:B:9:ASN:ND2	2.54	0.41
1:F:127:THR:HG22	1:F:127:THR:O	2.20	0.41
1:A:213:LEU:HA	1:A:214:PRO:HD3	1.81	0.41
1:C:99:HIS:O	1:C:101:PRO:HD3	2.21	0.41
1:C:175:MET:O	1:C:177:VAL:HG23	2.21	0.41
1:E:11:LYS:HA	1:E:11:LYS:HD3	1.72	0.41
1:E:153:VAL:HB	1:F:153:VAL:HG11	2.03	0.41
1:F:213:LEU:HA	1:F:214:PRO:HD3	1.74	0.41
1:E:203:LYS:HB3	1:E:203:LYS:HE2	1.93	0.40
1:F:28:LEU:HD23	1:F:31:PRO:HD3	2.04	0.40
1:B:181:GLU:H	1:B:181:GLU:CD	2.27	0.40
1:F:156:ARG:HG3	1:F:157:ALA:N	2.36	0.40
1:C:191:LEU:HD12	1:C:191:LEU:HA	1.95	0.40
1:D:191:LEU:O	1:D:194:ILE:HG22	2.22	0.40
1:F:20:CYS:SG	1:F:31:PRO:HB2	2.62	0.40
1:C:89:ASN:HB3	1:C:91:LEU:HD13	2.04	0.40
1:F:28:LEU:HD12	1:F:77:ARG:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	83/215 (39%)	81 (98%)	2 (2%)	0	100	100
1	B	206/215 (96%)	197 (96%)	9 (4%)	0	100	100
1	C	204/215 (95%)	194 (95%)	10 (5%)	0	100	100
1	D	22/215 (10%)	22 (100%)	0	0	100	100
1	E	161/215 (75%)	151 (94%)	10 (6%)	0	100	100
1	F	204/215 (95%)	192 (94%)	12 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	880/1290 (68%)	837 (95%)	43 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	58/190 (30%)	54 (93%)	4 (7%)	14	45
1	B	173/190 (91%)	167 (96%)	6 (4%)	32	64
1	C	24/190 (13%)	22 (92%)	2 (8%)	10	38
1	D	177/190 (93%)	167 (94%)	10 (6%)	19	52
1	E	172/190 (90%)	163 (95%)	9 (5%)	21	54
1	F	169/190 (89%)	154 (91%)	15 (9%)	9	34
All	All	773/1140 (68%)	727 (94%)	46 (6%)	17	50

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	159	THR
1	A	167	THR
1	A	191	LEU
1	A	211	LEU
1	B	95	GLU
1	B	106	ARG
1	B	133	ARG
1	B	153	VAL
1	B	159	THR
1	B	191	LEU
1	C	191	LEU
1	C	211	LEU
1	D	59	SER
1	D	67	LEU

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Mol	Chain	Res	Type
1	D	68	ARG
1	D	106	ARG
1	D	120	LEU
1	D	153	VAL
1	D	159	THR
1	D	177	VAL
1	D	191	LEU
1	D	211	LEU
1	E	9	ASN
1	E	68	ARG
1	E	72	GLN
1	E	91	LEU
1	E	121	PHE
1	E	159	THR
1	E	191	LEU
1	E	200	VAL
1	E	211	LEU
1	F	13	VAL
1	F	46	LEU
1	F	72	GLN
1	F	110	ARG
1	F	120	LEU
1	F	134	THR
1	F	153	VAL
1	F	159	THR
1	F	167	THR
1	F	168	ARG
1	F	170	VAL
1	F	185	GLN
1	F	194	ILE
1	F	210	LEU
1	F	211	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	89	ASN
1	B	199	ASN
1	D	112	HIS
1	E	9	ASN
1	E	116	HIS
1	E	185	GLN

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Mol	Chain	Res	Type
1	F	44	ASN
1	F	72	GLN
1	F	75	HIS
1	F	89	ASN
1	F	119	GLN
1	F	199	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	1T9	D	301	-	25,26,33	1.53	4 (16%)	28,31,38	1.35	3 (10%)
2	1T9	E	301	-	20,21,33	1.66	4 (20%)	23,26,38	1.48	2 (8%)
2	1T9	F	301	-	14,15,33	1.98	3 (21%)	17,20,38	1.65	2 (11%)
2	1T9	C	301	-	27,28,33	1.48	3 (11%)	30,33,38	1.34	3 (10%)
2	1T9	A	301	-	28,29,33	1.45	4 (14%)	31,34,38	1.32	3 (9%)
2	1T9	B	301	-	23,24,33	1.60	3 (13%)	26,29,38	1.39	3 (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	1T9	D	301	-	-	11/29/29/36	-
2	1T9	E	301	-	-	9/24/24/36	-
2	1T9	F	301	-	-	9/17/17/36	-
2	1T9	C	301	-	-	12/31/31/36	-
2	1T9	A	301	-	-	12/32/32/36	-
2	1T9	B	301	-	-	11/27/27/36	-

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	1T9	CAD-CAE	5.37	1.53	1.31
2	D	301	1T9	CAD-CAE	5.37	1.53	1.31
2	B	301	1T9	CAD-CAE	5.36	1.53	1.31
2	A	301	1T9	CAD-CAE	5.35	1.53	1.31
2	F	301	1T9	CAD-CAE	5.00	1.53	1.30
2	E	301	1T9	CAD-CAE	4.99	1.53	1.30
2	C	301	1T9	PAX-OAY	3.80	1.62	1.50
2	B	301	1T9	PAX-OAY	3.80	1.62	1.50
2	A	301	1T9	PAX-OAY	3.78	1.62	1.50
2	D	301	1T9	PAX-OAY	3.78	1.62	1.50
2	F	301	1T9	PAX-OAY	3.77	1.62	1.50
2	E	301	1T9	PAX-OAY	3.77	1.62	1.50
2	E	301	1T9	CAC-CAD	2.10	1.53	1.50
2	B	301	1T9	CAC-CAD	2.09	1.53	1.50
2	C	301	1T9	CAC-CAD	2.09	1.53	1.50
2	A	301	1T9	CAC-CAD	2.08	1.53	1.50
2	F	301	1T9	CAC-CAD	2.05	1.53	1.50
2	D	301	1T9	CAC-CAD	2.03	1.53	1.50
2	E	301	1T9	PAX-OAZ	2.02	1.62	1.54
2	D	301	1T9	PAX-OAZ	2.00	1.62	1.54
2	A	301	1T9	PAX-OAZ	2.00	1.62	1.54

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	301	1T9	CAC-CAD-CAE	-5.04	114.95	124.73
2	E	301	1T9	CAC-CAD-CAE	-5.01	115.01	124.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	1T9	CAC-CAD-CAE	-4.69	114.97	124.69
2	C	301	1T9	CAC-CAD-CAE	-4.69	114.98	124.69
2	A	301	1T9	CAC-CAD-CAE	-4.67	115.01	124.69
2	D	301	1T9	CAC-CAD-CAE	-4.64	115.07	124.69
2	F	301	1T9	OBA-PAX-OB	2.55	113.31	106.67
2	D	301	1T9	OBA-PAX-OB	2.54	113.29	106.67
2	C	301	1T9	OBA-PAX-OB	2.54	113.29	106.67
2	A	301	1T9	OBA-PAX-OB	2.54	113.28	106.67
2	E	301	1T9	OBA-PAX-OB	2.53	113.26	106.67
2	B	301	1T9	OBA-PAX-OB	2.52	113.25	106.67
2	B	301	1T9	CAQ-CAE-CAD	-2.36	114.97	125.47
2	C	301	1T9	CAQ-CAE-CAD	-2.35	114.99	125.47
2	A	301	1T9	CAQ-CAE-CAD	-2.35	115.01	125.47
2	D	301	1T9	CAQ-CAE-CAD	-2.35	115.01	125.47

There are no chirality outliers.

All (64) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	1T9	CAA-OB-PAX-OAY
2	A	301	1T9	CAA-OB-PAX-OAZ
2	A	301	1T9	CAA-OB-PAX-OBA
2	B	301	1T9	CAC-CAB-NAW-CAU
2	B	301	1T9	CAA-OB-PAX-OAY
2	B	301	1T9	CAA-OB-PAX-OAZ
2	B	301	1T9	CAA-OB-PAX-OBA
2	C	301	1T9	CAA-OB-PAX-OAZ
2	C	301	1T9	CAA-OB-PAX-OBA
2	D	301	1T9	CAA-OB-PAX-OAY
2	D	301	1T9	CAA-OB-PAX-OAZ
2	D	301	1T9	CAA-OB-PAX-OBA
2	E	301	1T9	CAA-OB-PAX-OAY
2	E	301	1T9	CAA-OB-PAX-OAZ
2	E	301	1T9	CAA-OB-PAX-OBA
2	F	301	1T9	CAA-OB-PAX-OAY
2	F	301	1T9	CAA-OB-PAX-OAZ
2	F	301	1T9	CAA-OB-PAX-OBA
2	B	301	1T9	CAU-CAT-CBC-CBD
2	A	301	1T9	CAU-CAT-CBC-CBD
2	D	301	1T9	CAM-CAN-CAO-CAP
2	B	301	1T9	CAA-CAB-NAW-CAU
2	A	301	1T9	CAM-CAN-CAO-CAP

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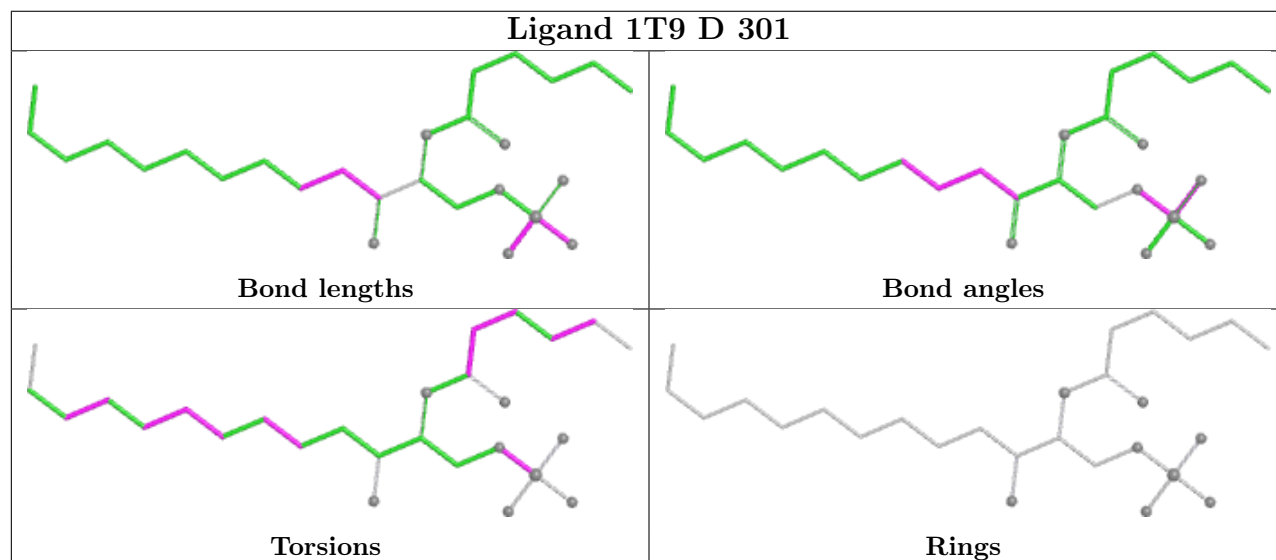
Mol	Chain	Res	Type	Atoms
2	A	301	1T9	CBC-CBD-CBE-CBF
2	E	301	1T9	CBD-CBE-CBF-CBG
2	C	301	1T9	CBD-CBE-CBF-CBG
2	C	301	1T9	CAN-CAO-CAP-CAQ
2	A	301	1T9	CBD-CBE-CBF-CBG
2	B	301	1T9	CAL-CAM-CAN-CAO
2	C	301	1T9	OAR-CAC-CAD-CAE
2	F	301	1T9	OAR-CAC-CAD-CAE
2	C	301	1T9	CAK-CAL-CAM-CAN
2	C	301	1T9	CAA-OBB-PAX-OAY
2	C	301	1T9	CAJ-CAK-CAL-CAM
2	C	301	1T9	CBC-CBD-CBE-CBF
2	D	301	1T9	CAU-CAT-CBC-CBD
2	D	301	1T9	CAN-CAO-CAP-CAQ
2	A	301	1T9	CBE-CBF-CBG-CBH
2	B	301	1T9	CAK-CAL-CAM-CAN
2	A	301	1T9	OBB-CAA-CAB-NAW
2	B	301	1T9	OBB-CAA-CAB-NAW
2	E	301	1T9	OBB-CAA-CAB-NAW
2	F	301	1T9	OBB-CAA-CAB-NAW
2	F	301	1T9	CAB-CAA-OBB-PAX
2	D	301	1T9	CAK-CAL-CAM-CAN
2	C	301	1T9	CBC-CAT-CAU-OAV
2	A	301	1T9	OBB-CAA-CAB-CAC
2	B	301	1T9	OBB-CAA-CAB-CAC
2	E	301	1T9	OBB-CAA-CAB-CAC
2	E	301	1T9	CAT-CBC-CBD-CBE
2	F	301	1T9	CAB-CAC-CAD-CAE
2	F	301	1T9	CAA-CAB-NAW-CAU
2	C	301	1T9	CBC-CAT-CAU-NAW
2	A	301	1T9	CAI-CAJ-CAK-CAL
2	D	301	1T9	CBC-CAT-CAU-OAV
2	D	301	1T9	CAD-CAE-CAQ-CAP
2	C	301	1T9	CAO-CAP-CAQ-CAE
2	D	301	1T9	CBC-CAT-CAU-NAW
2	A	301	1T9	CAD-CAE-CAQ-CAP
2	E	301	1T9	CBC-CAT-CAU-OAV
2	E	301	1T9	CBC-CAT-CAU-NAW
2	F	301	1T9	OBB-CAA-CAB-CAC
2	B	301	1T9	CAM-CAN-CAO-CAP
2	D	301	1T9	CBC-CBD-CBE-CBF

There are no ring outliers.

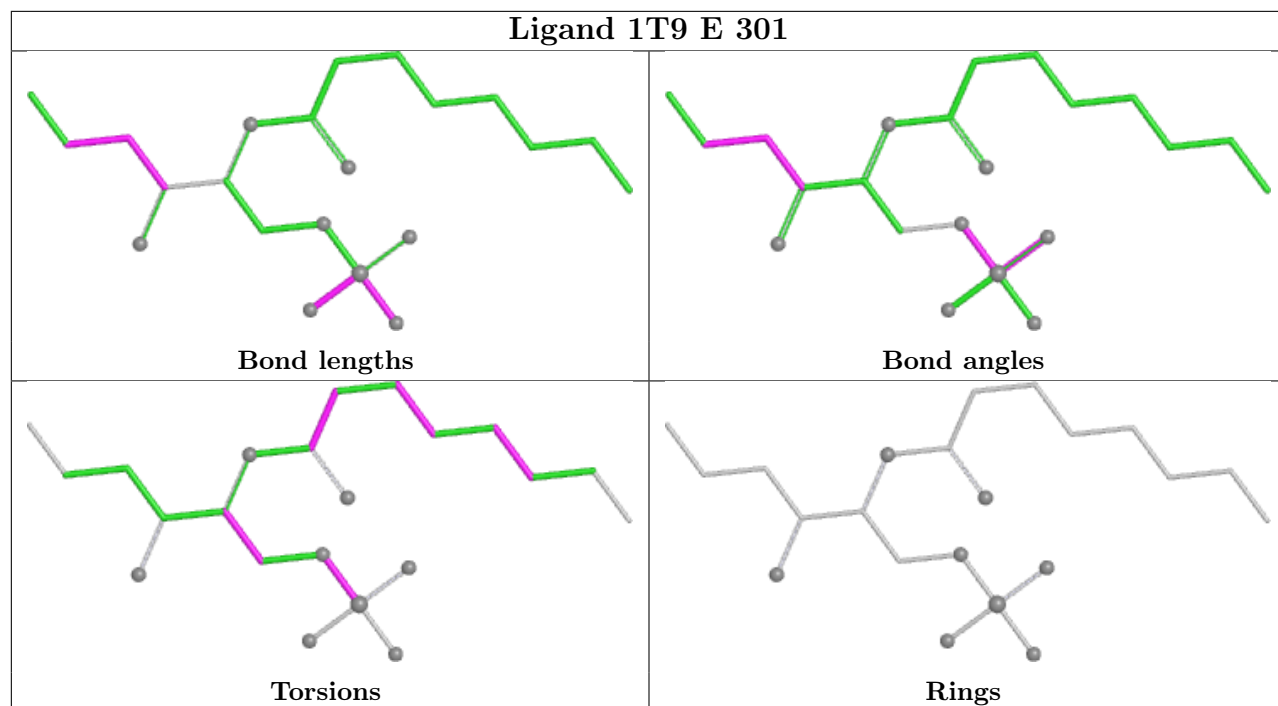
5 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	301	1T9	2	0
2	E	301	1T9	1	0
2	F	301	1T9	2	0
2	C	301	1T9	6	0
2	A	301	1T9	4	0

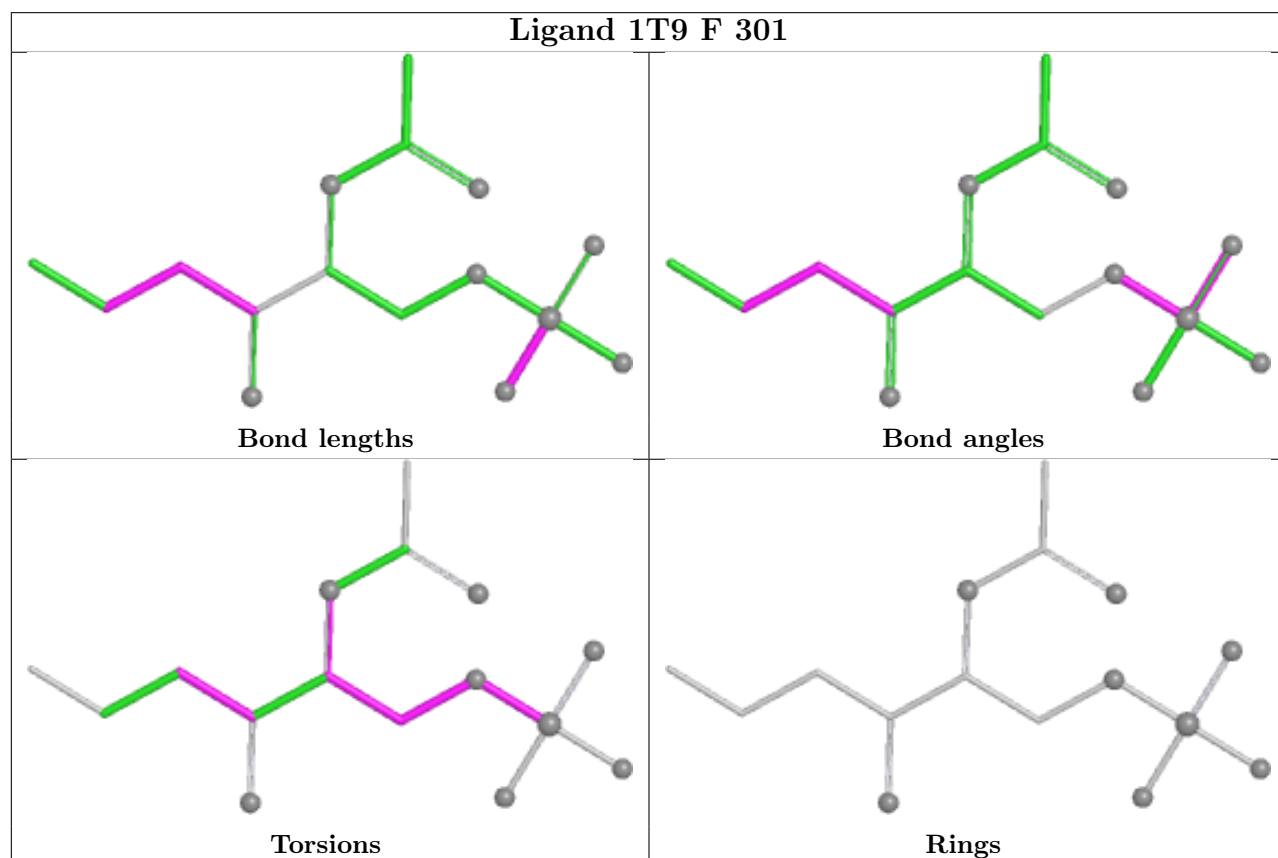
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

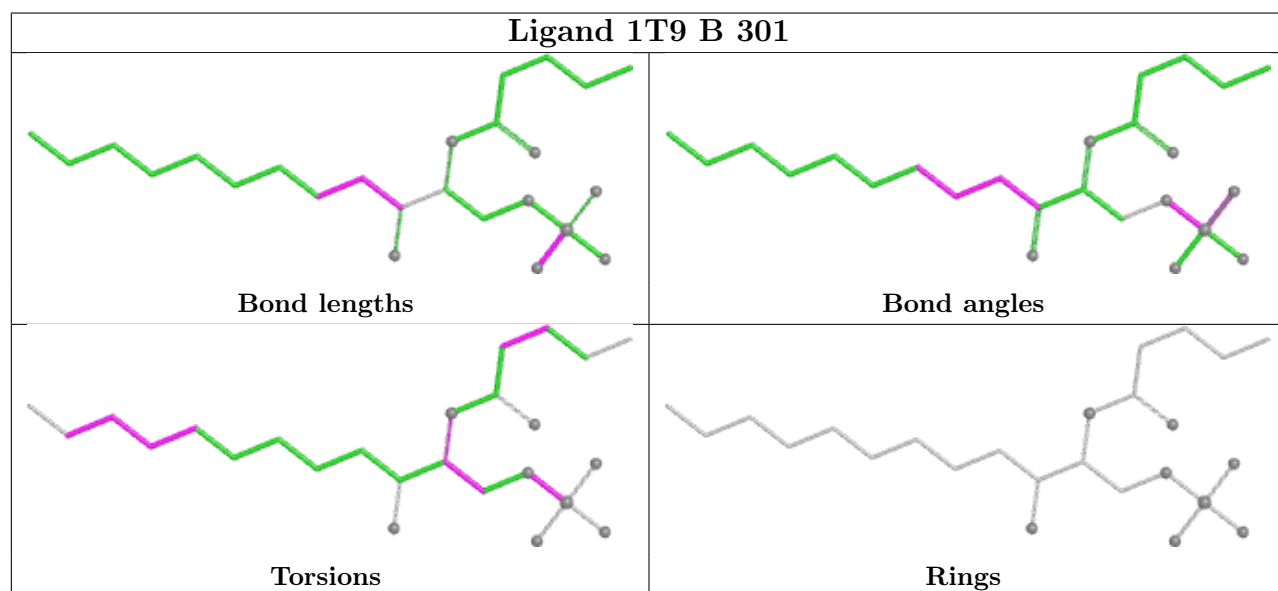
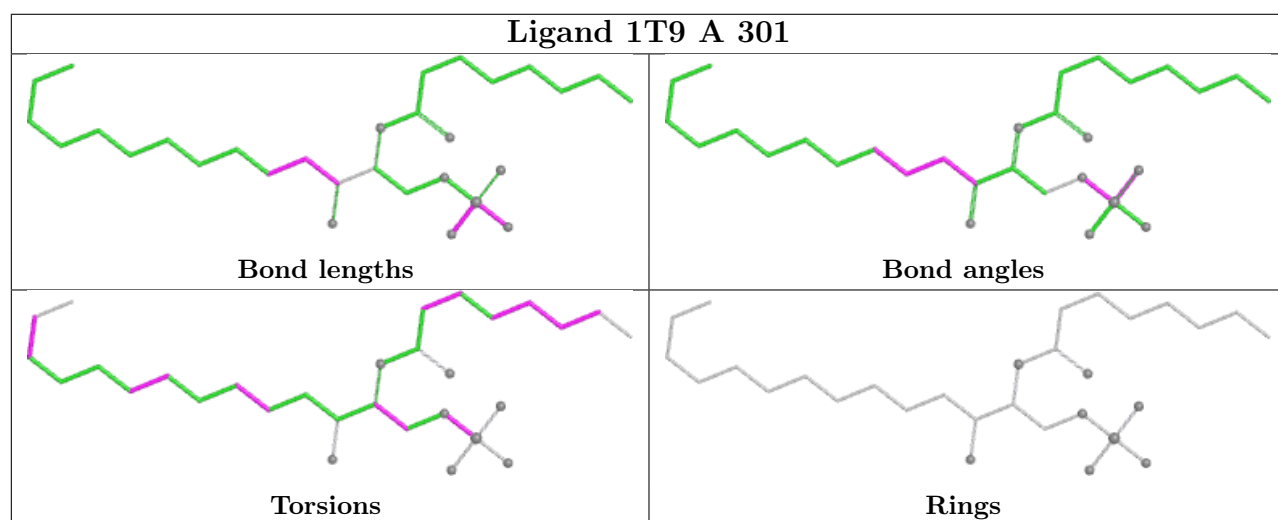
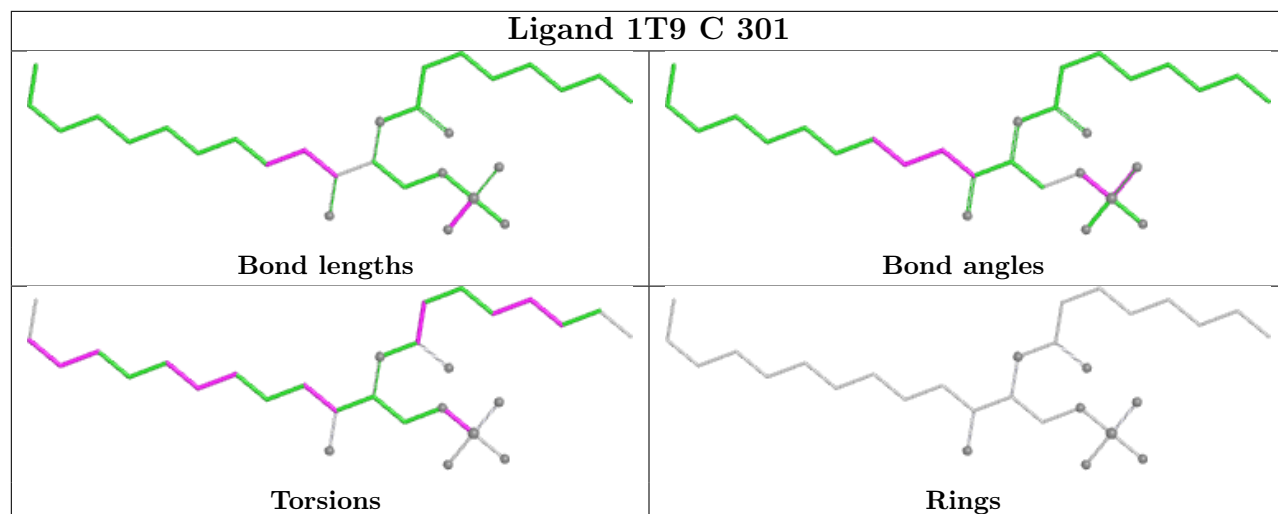


Ligand 1T9 E 301



Ligand 1T9 F 301





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	206/215 (95%)	-0.49	0	100 100	28, 54, 98, 117	0
1	B	208/215 (96%)	-0.44	0	100 100	30, 63, 107, 157	0
1	C	206/215 (95%)	-0.50	0	100 100	32, 62, 108, 147	0
1	D	206/215 (95%)	-0.46	0	100 100	31, 72, 132, 202	0
1	E	206/215 (95%)	-0.19	1 (0%)	87 76	41, 79, 130, 168	0
1	F	206/215 (95%)	0.04	0	100 100	57, 136, 195, 236	0
All	All	1238/1290 (95%)	-0.34	1 (0%)	92 89	28, 72, 161, 236	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	43	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

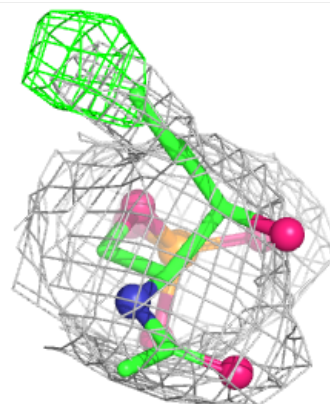
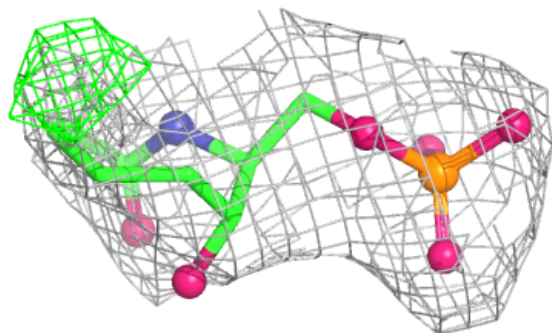
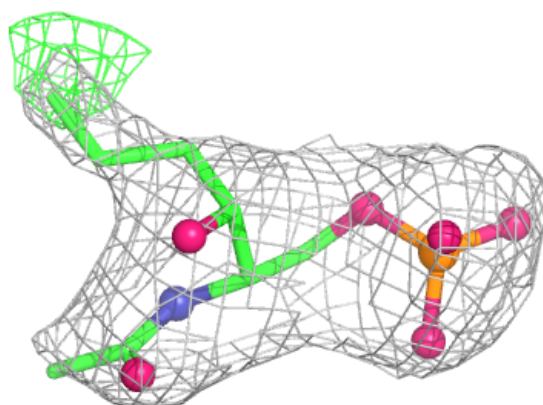
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	1T9	F	301	16/34	0.80	0.11	89,94,96,98	0
2	1T9	B	301	25/34	0.92	0.11	54,57,61,61	0
2	1T9	C	301	29/34	0.93	0.12	54,55,59,60	0
2	1T9	D	301	27/34	0.94	0.10	40,43,47,47	0
2	1T9	E	301	22/34	0.94	0.09	54,57,61,62	0
2	1T9	A	301	30/34	0.94	0.10	34,35,37,38	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.

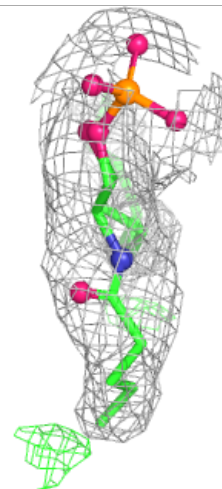
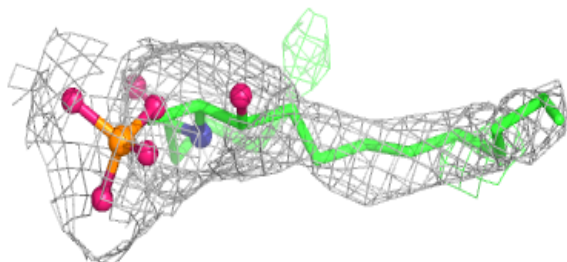
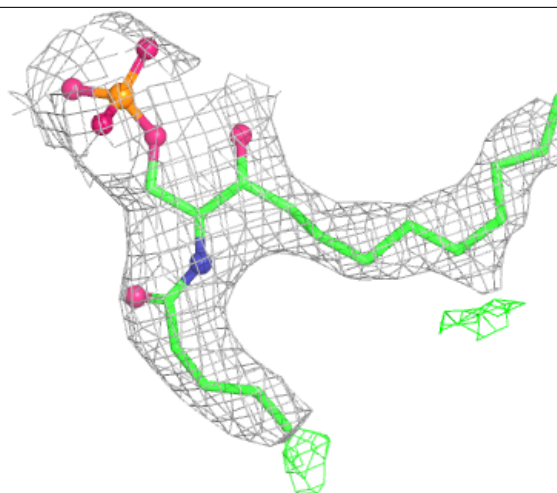
Electron density around 1T9 F 301:

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



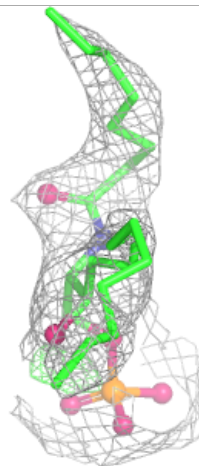
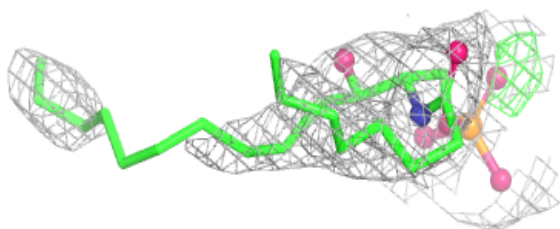
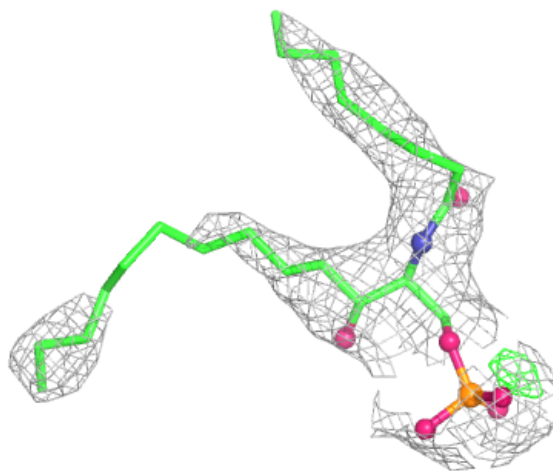
Electron density around 1T9 B 301:

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



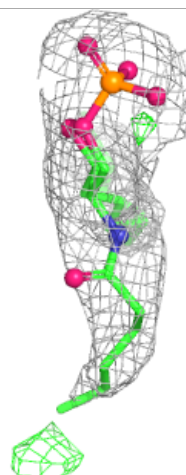
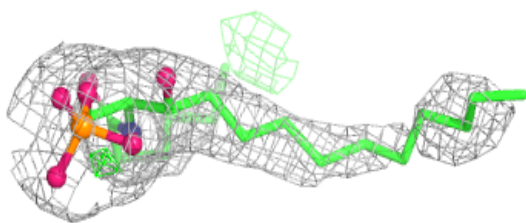
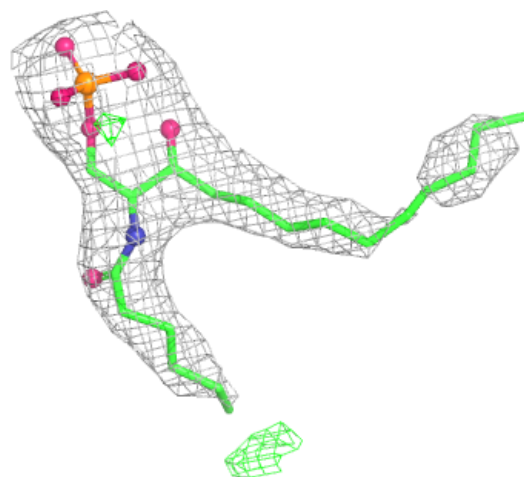
Electron density around 1T9 C 301:

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



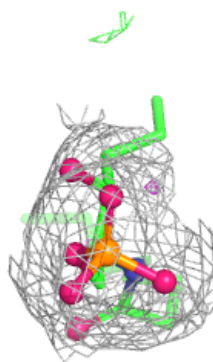
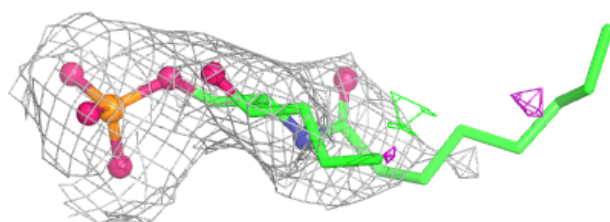
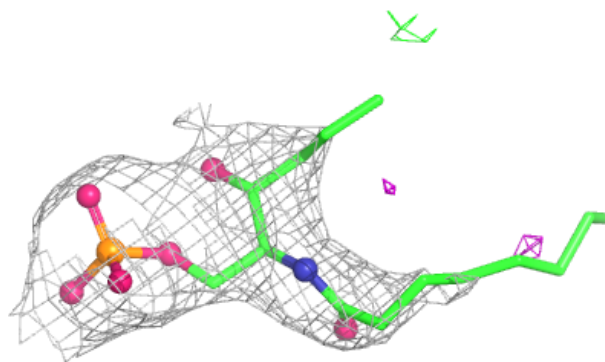
Electron density around 1T9 D 301:

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



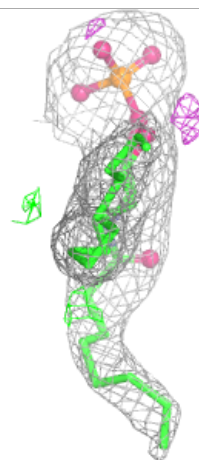
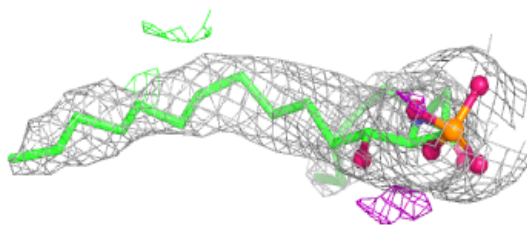
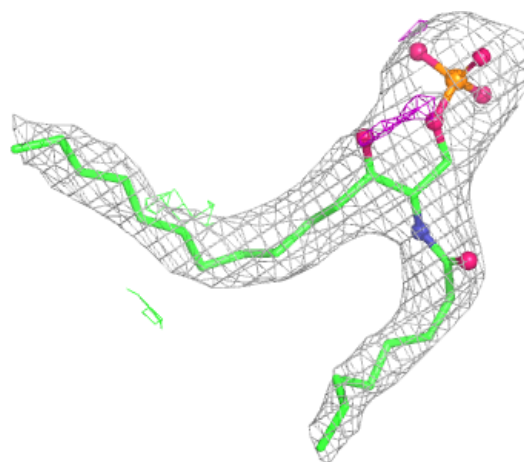
Electron density around 1T9 E 301:

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



Electron density around 1T9 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.