



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 03:15 AM UTC

PDB ID : 9I55 / pdb_00009i55
Title : Mouse phosphomannomutase 2 in complex with the activator glucose-1,6-bisphosphate
Authors : Del Cano-Ochoa, F.; Vilar, M.; Vilas, A.; Company, R.; Perez, B.; Ramon-Maiques, S.
Deposited on : 2025-01-27
Resolution : 2.75 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

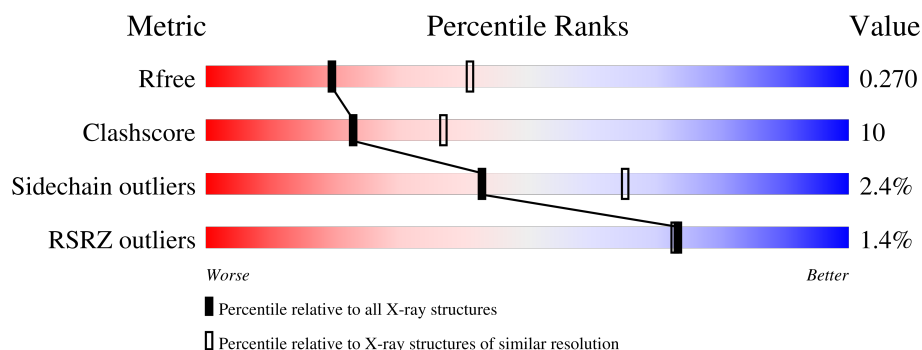
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	244	<div> <div>88%</div> <div>11% .</div> </div>
1	B	244	<div> <div>85%</div> <div>15%</div> </div>
1	C	244	<div> <div>82%</div> <div>18%</div> </div>
1	D	244	<div> <div>%</div> <div>73%</div> <div>26% .</div> </div>
1	E	244	<div> <div>3%</div> <div>70%</div> <div>28% ..</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	244	<div><div></div><div>5%</div><div>68%</div><div>29%</div><div>..</div></div>

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 23185 atoms, of which 11285 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 Phosphomannomutase 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	242	Total	C	H	N	O	S	0	0	0
			3839	1232	1904	330	363	10			
1	B	243	Total	C	H	N	O	S	0	0	0
			3878	1242	1929	332	364	11			
1	D	242	Total	C	H	N	O	S	0	0	0
			3799	1224	1880	324	361	10			
1	F	241	Total	C	H	N	O	S	0	0	0
			3631	1187	1770	315	349	10			
1	C	243	Total	C	H	N	O	S	0	0	0
			3863	1239	1918	331	364	11			
1	E	242	Total	C	H	N	O	S	0	0	0
			3730	1209	1836	325	350	10			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q9Z2M7
A	0	PRO	-	expression tag	UNP Q9Z2M7
B	-1	GLY	-	expression tag	UNP Q9Z2M7
B	0	PRO	-	expression tag	UNP Q9Z2M7
D	-1	GLY	-	expression tag	UNP Q9Z2M7
D	0	PRO	-	expression tag	UNP Q9Z2M7
F	-1	GLY	-	expression tag	UNP Q9Z2M7
F	0	PRO	-	expression tag	UNP Q9Z2M7
C	-1	GLY	-	expression tag	UNP Q9Z2M7
C	0	PRO	-	expression tag	UNP Q9Z2M7
E	-1	GLY	-	expression tag	UNP Q9Z2M7
E	0	PRO	-	expression tag	UNP Q9Z2M7

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Mg	0	0
			3	3		
3	B	1	Total	Mg	0	0
			1	1		
3	D	2	Total	Mg	0	0
			2	2		
3	F	2	Total	Mg	0	0
			2	2		
3	C	2	Total	Mg	0	0
			2	2		
3	E	2	Total	Mg	0	0
			2	2		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		
4	B	1	Total	Cl	0	0
			1	1		

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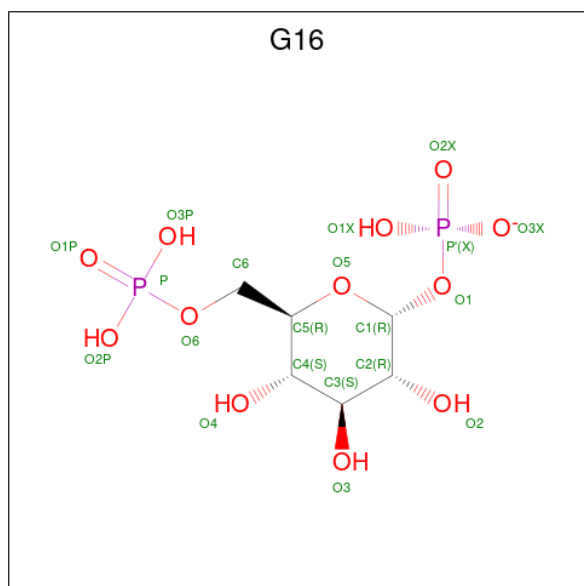
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	E	2	Total	Cl			0	0
			2	2				

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	Na			0	0
			1	1				

- Molecule 6 is 1,6-di-O-phosphono-alpha-D-glucopyranose (CCD ID: G16) (formula: C₆H₁₃O₁₂P₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	H	O	P	0	0
			30	6	10	12	2		
6	D	1	Total	C	H	O	P	0	0
			30	6	10	12	2		
6	F	1	Total	C	H	O	P	0	0
			30	6	10	12	2		
6	C	1	Total	C	H	O	P	0	0
			30	6	10	12	2		

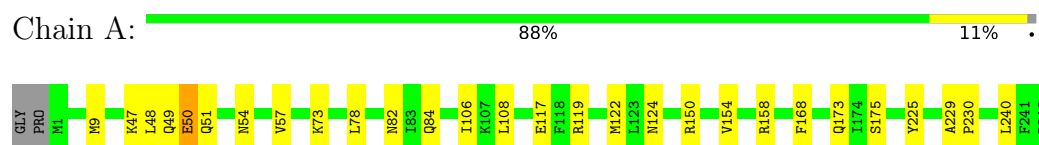
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	73	Total 73	O 73	0	0
7	B	59	Total 59	O 59	0	0
7	D	52	Total 52	O 52	0	0
7	F	20	Total 20	O 20	0	0
7	C	60	Total 60	O 60	0	0
7	E	30	Total 30	O 30	0	0

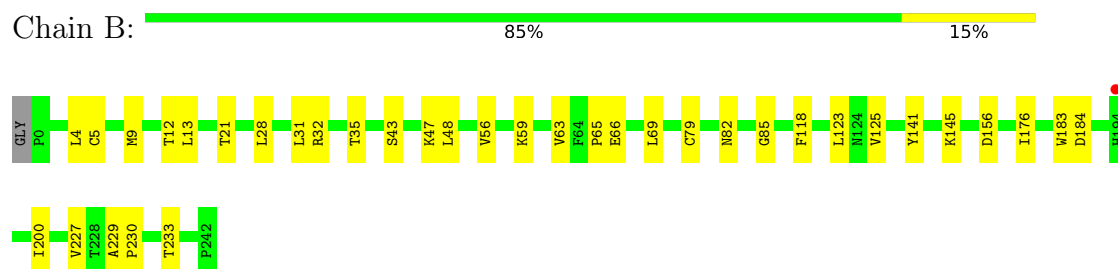
3 Residue-property plots

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

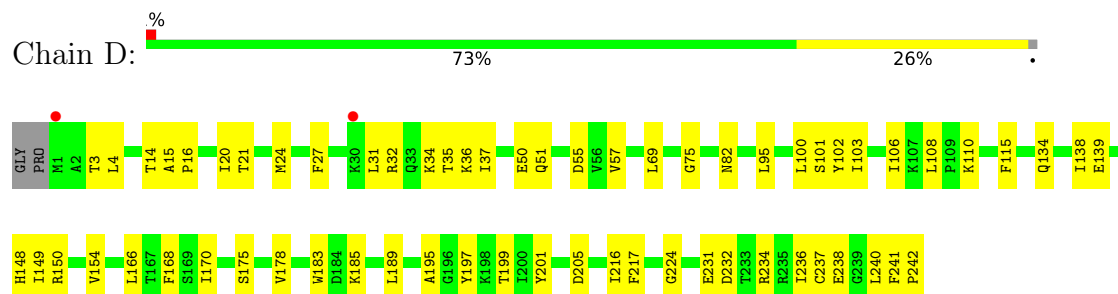
• Molecule 1: Phosphomannomutase 2



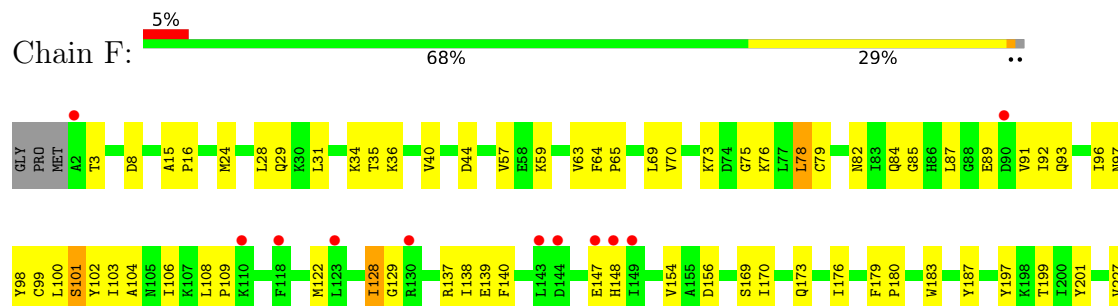
• Molecule 1: Phosphomannomutase 2



• Molecule 1: Phosphomannomutase 2



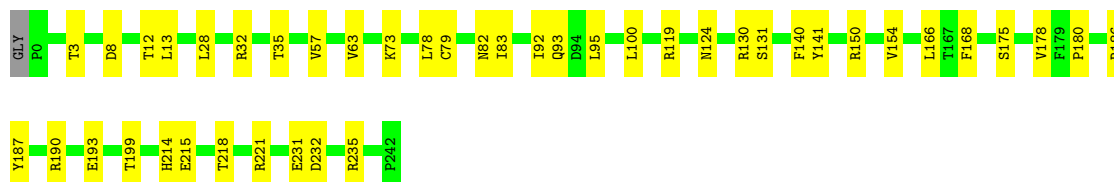
• Molecule 1: Phosphomannomutase 2





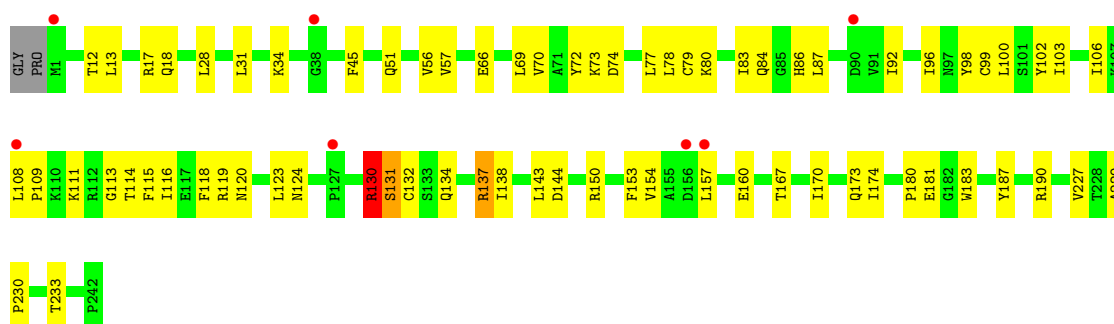
● Molecule 1: Phosphomannomutase 2

Chain C: 82% 18%



● Molecule 1: Phosphomannomutase 2

Chain E: 3% 70% 28% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.28Å 99.14Å 213.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	64.67 – 2.75 64.67 – 2.75	Depositor EDS
% Data completeness (in resolution range)	98.8 (64.67-2.75) 98.8 (64.67-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.19 (at 2.73Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.205 , 0.265 0.217 , 0.270	Depositor DCC
R_{free} test set	2289 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	61.4	Xtriage
Anisotropy	0.377	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	23185	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% 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: GOL, MG, G16, CL, NA

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	0.20	0/1975	0.39	0/2656
1	B	0.19	0/1990	0.37	0/2674
1	C	0.20	0/1986	0.36	0/2670
1	D	0.19	0/1959	0.39	0/2637
1	E	0.25	0/1934	0.42	0/2608
1	F	0.20	0/1900	0.34	0/2568
All	All	0.21	0/11744	0.38	0/15813

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	B	0	1
1	E	0	3
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	32	ARG	Sidechain
1	E	130	ARG	Sidechain
1	E	137	ARG	Sidechain
1	E	17	ARG	Sidechain

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	1935	1904	1906	17	0
1	B	1949	1929	1931	21	0
1	C	1945	1918	1920	23	0
1	D	1919	1880	1880	55	1
1	E	1894	1836	1838	64	0
1	F	1861	1770	1768	60	0
2	A	6	8	8	0	0
3	A	3	0	0	0	0
3	B	1	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
3	E	2	0	0	0	0
3	F	2	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	E	2	0	0	2	0
5	A	1	0	0	0	1
6	B	20	10	10	0	0
6	C	20	10	10	1	0
6	D	20	10	10	0	0
6	F	20	10	10	1	0
7	A	73	0	0	1	0
7	B	59	0	0	0	0
7	C	60	0	0	0	0
7	D	52	0	0	2	0
7	E	30	0	0	1	0
7	F	20	0	0	1	0
All	All	11900	11285	11291	238	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 10.

All (238) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:236:ILE:HG22	1:D:240:LEU:HD11	1.55	0.85
1:D:150:ARG:O	1:D:154:VAL:HG23	1.85	0.77
1:C:231:GLU:OE1	1:C:231:GLU:N	2.18	0.76
1:E:153:PHE:CE1	1:E:157:LEU:HD11	2.21	0.75
1:B:48:LEU:HD12	1:B:65:PRO:HB3	1.71	0.73
1:D:31:LEU:CD1	1:D:35:THR:HG22	2.21	0.71
1:A:47:LYS:NZ	1:A:50:GLU:OE2	2.24	0.70
1:E:13:LEU:HD22	1:E:28:LEU:HD11	1.74	0.69
4:E:304:CL:CL	7:E:417:HOH:O	2.47	0.69
1:C:13:LEU:HD22	1:C:28:LEU:HD11	1.74	0.69
1:F:8:ASP:OD2	7:F:401:HOH:O	2.11	0.68
1:F:96:ILE:HG22	1:F:100:LEU:HD11	1.76	0.68
1:B:9:MET:SD	1:B:48:LEU:HD23	2.35	0.66
1:D:240:LEU:HD12	1:D:240:LEU:H	1.58	0.66
1:E:83:ILE:O	1:E:87:LEU:HD23	1.96	0.66
1:A:57:VAL:HG22	7:A:407:HOH:O	1.95	0.66
1:B:13:LEU:HD22	1:B:28:LEU:HD11	1.78	0.66
1:B:63:VAL:HG12	1:B:65:PRO:HD3	1.79	0.65
1:D:31:LEU:HD11	1:D:35:THR:HG22	1.79	0.65
1:E:83:ILE:HG13	1:E:87:LEU:HD23	1.81	0.63
1:F:97:ASN:HA	1:F:100:LEU:HD12	1.81	0.63
1:F:102:TYR:CZ	1:F:106:ILE:HD11	2.35	0.62
1:D:236:ILE:O	1:D:240:LEU:HD12	1.99	0.61
1:F:31:LEU:HD11	1:F:35:THR:HG22	1.82	0.61
1:D:3:THR:HG23	1:D:35:THR:OG1	1.99	0.61
1:A:119:ARG:NH1	1:A:124:ASN:OD1	2.33	0.61
1:F:96:ILE:CG2	1:F:100:LEU:HD11	2.30	0.61
1:D:24:MET:HE2	1:D:24:MET:HA	1.82	0.60
1:C:168:PHE:CD2	1:C:178:VAL:HG22	2.37	0.60
1:E:87:LEU:N	1:E:87:LEU:HD22	2.17	0.59
1:E:113:GLY:O	1:E:114:THR:OG1	2.16	0.59
1:D:15:ALA:HB1	1:D:16:PRO:HD2	1.85	0.59
1:F:108:LEU:HG	1:F:109:PRO:CD	2.34	0.58
1:E:150:ARG:NH1	1:E:173:GLN:O	2.37	0.58
1:F:96:ILE:HG22	1:F:100:LEU:CD1	2.35	0.57
1:D:236:ILE:HD12	1:D:236:ILE:H	1.68	0.57
1:B:12:THR:HG23	1:B:227:VAL:HG21	1.87	0.56
1:C:150:ARG:O	1:C:154:VAL:HG23	2.05	0.56
1:F:108:LEU:HG	1:F:109:PRO:HD3	1.87	0.56
1:F:69:LEU:HD21	1:F:183:TRP:CE3	2.40	0.56
1:F:69:LEU:HD21	1:F:183:TRP:CZ3	2.39	0.56
1:D:31:LEU:HD11	1:D:35:THR:CG2	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:82:ASN:OD1	1:F:85:GLY:N	2.38	0.56
1:D:3:THR:HG23	1:D:35:THR:HG1	1.68	0.55
1:A:73:LYS:HB3	1:A:78:LEU:HD13	1.89	0.55
1:E:132:CYS:SG	1:E:137:ARG:HG3	2.47	0.55
1:E:73:LYS:NZ	1:E:74:ASP:OD2	2.39	0.55
1:A:9:MET:HE1	1:A:48:LEU:HA	1.89	0.55
1:A:122:MET:HE2	1:A:124:ASN:ND2	2.22	0.55
1:D:189:LEU:HD11	1:D:216:ILE:HD12	1.89	0.55
1:F:99:CYS:O	1:F:103:ILE:HD12	2.06	0.55
1:E:18:GLN:O	1:E:51:GLN:NE2	2.40	0.54
1:D:240:LEU:HD12	1:D:240:LEU:N	2.23	0.54
1:E:70:VAL:HG22	1:E:80:LYS:CG	2.38	0.54
1:D:15:ALA:HB1	1:D:16:PRO:CD	2.37	0.54
1:E:134:GLN:O	1:E:138:ILE:HG12	2.06	0.54
1:B:43:SER:HB3	1:B:47:LYS:HB3	1.89	0.54
1:B:118:PHE:CD2	1:B:123:LEU:HD21	2.42	0.54
1:B:66:GLU:O	1:B:69:LEU:HD12	2.08	0.54
1:F:236:ILE:O	1:F:240:LEU:HD12	2.07	0.53
1:A:49:GLN:HG2	1:A:54:ASN:HA	1.91	0.53
1:F:147:GLU:O	1:F:148:HIS:CB	2.56	0.53
1:F:102:TYR:O	1:F:106:ILE:HG12	2.09	0.52
1:D:50:GLU:OE1	1:D:51:GLN:HG3	2.10	0.52
1:C:214:HIS:O	1:C:218:THR:HG23	2.09	0.52
1:E:109:PRO:HG3	1:E:143:LEU:HD12	1.91	0.52
1:D:154:VAL:HG21	1:D:170:ILE:HG13	1.91	0.52
1:F:69:LEU:CD2	1:F:183:TRP:CZ3	2.93	0.51
1:C:119:ARG:NH1	1:C:124:ASN:OD1	2.43	0.51
1:B:69:LEU:HD21	1:B:183:TRP:CE3	2.45	0.51
1:D:69:LEU:HD11	1:D:183:TRP:CE3	2.46	0.51
1:B:4:LEU:HD23	1:B:200:ILE:HG12	1.91	0.51
1:E:115:PHE:C	1:E:116:ILE:HG13	2.36	0.51
1:F:128:ILE:HG23	1:F:129:GLY:N	2.26	0.51
1:E:167:THR:OG1	1:E:181:GLU:HG2	2.11	0.51
1:E:108:LEU:HB3	1:E:109:PRO:HD2	1.93	0.50
1:E:173:GLN:HG2	1:E:174:ILE:HG23	1.92	0.50
1:E:153:PHE:CD1	1:E:157:LEU:HD11	2.46	0.50
1:E:102:TYR:O	1:E:106:ILE:HD12	2.12	0.50
1:D:205:ASP:OD1	7:D:401:HOH:O	2.20	0.50
1:F:106:ILE:O	1:F:108:LEU:N	2.40	0.49
1:E:115:PHE:N	4:E:303:CL:CL	2.82	0.49
1:F:98:TYR:CD1	1:F:98:TYR:C	2.90	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:102:TYR:CD1	1:E:106:ILE:HD11	2.46	0.49
1:A:225:TYR:CE2	1:A:240:LEU:HD11	2.48	0.49
1:A:150:ARG:O	1:A:154:VAL:HG23	2.13	0.49
1:B:5:CYS:SG	1:B:35:THR:HG21	2.52	0.49
1:A:106:ILE:HG22	1:A:108:LEU:HG	1.94	0.49
1:F:108:LEU:HD21	1:F:140:PHE:CE1	2.48	0.49
1:F:232:ASP:OD1	1:F:235:ARG:NH2	2.45	0.49
1:C:186:ARG:HE	1:C:215:GLU:HB3	1.77	0.49
1:F:101:SER:O	1:F:104:ALA:HB3	2.13	0.49
1:D:168:PHE:CD2	1:D:178:VAL:HG22	2.48	0.49
1:F:180:PRO:HB2	1:F:183:TRP:CD1	2.47	0.49
1:E:83:ILE:HG13	1:E:87:LEU:CD2	2.43	0.48
1:D:57:VAL:HG22	1:D:75:GLY:HA2	1.95	0.48
1:F:89:GLU:HA	1:F:92:ILE:HB	1.94	0.48
1:E:153:PHE:CE1	1:E:157:LEU:CD1	2.95	0.48
1:D:31:LEU:HD12	1:D:31:LEU:O	2.14	0.48
1:D:102:TYR:HE1	1:D:106:ILE:HD11	1.78	0.48
1:D:110:LYS:NZ	1:D:139:GLU:OE1	2.46	0.48
1:F:31:LEU:CD1	1:F:35:THR:HG22	2.43	0.47
1:E:72:TYR:CD1	1:E:77:LEU:HA	2.49	0.47
1:E:92:ILE:O	1:E:96:ILE:HG13	2.14	0.47
1:E:144:ASP:OD2	1:E:150:ARG:HG3	2.14	0.47
1:D:236:ILE:CG2	1:D:240:LEU:HD11	2.37	0.47
1:E:12:THR:O	1:E:230:PRO:HA	2.14	0.47
1:E:56:VAL:HG13	1:E:57:VAL:N	2.29	0.47
1:A:47:LYS:HE3	1:A:51:GLN:OE1	2.13	0.47
1:E:102:TYR:CE1	1:E:106:ILE:HD11	2.50	0.47
1:E:73:LYS:HB3	1:E:78:LEU:HD12	1.96	0.47
1:F:57:VAL:HG12	1:F:75:GLY:HA2	1.98	0.46
1:C:83:ILE:HG12	1:C:92:ILE:CD1	2.46	0.46
1:D:32:ARG:HD2	1:D:37:ILE:HD11	1.97	0.46
1:F:31:LEU:O	1:F:34:LYS:N	2.49	0.46
1:A:47:LYS:HZ3	1:A:50:GLU:CD	2.21	0.46
1:D:24:MET:HA	1:D:24:MET:CE	2.45	0.46
1:C:83:ILE:HG12	1:C:92:ILE:HD11	1.97	0.46
1:F:29:GLN:NE2	1:F:59:LYS:HG2	2.31	0.46
1:A:229:ALA:HB1	1:A:230:PRO:HD2	1.97	0.46
1:E:77:LEU:HD23	1:E:79:CYS:N	2.31	0.46
1:A:158:ARG:NH2	1:A:168:PHE:O	2.49	0.46
1:D:4:LEU:HD12	1:D:36:LYS:O	2.16	0.46
1:D:102:TYR:CE1	1:D:106:ILE:HD11	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:189:LEU:HD21	1:D:216:ILE:HD11	1.98	0.45
1:C:190:ARG:HG3	1:C:190:ARG:HH11	1.81	0.45
1:C:193:GLU:OE2	1:C:221:ARG:NH1	2.49	0.45
1:D:195:ALA:HB1	1:D:197:TYR:HE1	1.82	0.45
1:F:100:LEU:HD21	1:E:115:PHE:CD1	2.51	0.45
1:B:229:ALA:HB1	1:B:230:PRO:HD2	1.98	0.45
1:B:200:ILE:O	1:B:200:ILE:HG22	2.17	0.45
1:E:31:LEU:HA	1:E:34:LYS:HE3	1.99	0.45
1:E:66:GLU:OE2	1:E:120:ASN:ND2	2.49	0.45
1:E:229:ALA:HB1	1:E:230:PRO:HD2	1.99	0.45
1:E:12:THR:HG23	1:E:227:VAL:HG21	1.98	0.45
1:D:232:ASP:O	1:D:236:ILE:HD12	2.17	0.45
1:F:63:VAL:HG12	1:F:65:PRO:HD3	1.99	0.44
1:D:100:LEU:HD22	1:C:100:LEU:HD23	1.98	0.44
1:D:234:ARG:O	1:D:238:GLU:HG2	2.17	0.44
1:F:73:LYS:HB2	1:F:78:LEU:HD12	1.99	0.44
1:F:93:GLN:HG2	1:E:111:LYS:HB2	1.98	0.44
1:F:154:VAL:HG22	1:F:176:ILE:HG21	1.99	0.44
1:C:32:ARG:HA	1:C:35:THR:O	2.17	0.44
1:A:229:ALA:HB1	1:A:230:PRO:CD	2.47	0.44
1:F:87:LEU:HD23	1:F:92:ILE:CD1	2.46	0.44
1:F:87:LEU:HG	1:F:91:VAL:HG13	1.98	0.44
1:E:227:VAL:HG11	1:E:233:THR:HG23	2.00	0.44
1:D:50:GLU:OE1	1:D:51:GLN:CG	2.66	0.44
1:E:98:TYR:CD1	1:E:98:TYR:C	2.95	0.44
1:D:103:ILE:HD13	1:D:115:PHE:HB3	1.99	0.44
1:E:99:CYS:O	1:E:102:TYR:N	2.50	0.44
1:B:31:LEU:HD12	1:B:31:LEU:O	2.17	0.44
1:E:70:VAL:HG22	1:E:80:LYS:HG3	2.00	0.44
1:E:99:CYS:O	1:E:100:LEU:C	2.61	0.44
1:B:118:PHE:HA	1:B:123:LEU:HD23	2.00	0.43
1:E:100:LEU:HG	1:E:116:ILE:CD1	2.48	0.43
1:F:82:ASN:ND2	1:F:84:GLN:NE2	2.66	0.43
1:B:229:ALA:HB1	1:B:230:PRO:CD	2.48	0.43
1:D:236:ILE:HD12	1:D:236:ILE:N	2.33	0.43
1:E:87:LEU:HD21	1:E:180:PRO:HG3	2.00	0.43
1:E:118:PHE:CE2	1:E:123:LEU:HD21	2.53	0.43
1:A:117:GLU:HB3	1:A:124:ASN:HB2	2.01	0.43
1:B:69:LEU:HD21	1:B:183:TRP:CZ3	2.54	0.43
1:D:185:LYS:NZ	7:D:405:HOH:O	2.52	0.43
1:F:29:GLN:OE1	1:F:29:GLN:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:130:ARG:NH2	6:C:301:G16:O1	2.51	0.43
1:E:87:LEU:N	1:E:87:LEU:CD2	2.82	0.43
1:C:140:PHE:O	1:C:141:TYR:C	2.61	0.43
1:E:69:LEU:HD21	1:E:183:TRP:CE3	2.53	0.43
1:E:84:GLN:OE1	1:E:84:GLN:N	2.51	0.43
1:F:3:THR:HG1	1:F:35:THR:HG1	1.66	0.43
1:F:137:ARG:CD	1:F:173:GLN:OE1	2.67	0.43
1:C:57:VAL:HA	1:C:63:VAL:HG21	2.00	0.43
1:E:100:LEU:HG	1:E:116:ILE:HD11	2.01	0.43
1:E:144:ASP:OD2	1:E:150:ARG:HB2	2.19	0.43
1:C:95:LEU:HD11	1:C:166:LEU:HD13	2.00	0.43
1:D:217:PHE:CD2	1:D:224:GLY:HA3	2.53	0.42
1:D:199:THR:HG21	1:D:201:TYR:CE1	2.55	0.42
1:F:44:ASP:OD1	1:F:44:ASP:N	2.52	0.42
1:E:154:VAL:HG21	1:E:170:ILE:HG13	2.01	0.42
1:D:236:ILE:O	1:D:240:LEU:CD1	2.67	0.42
1:E:113:GLY:C	1:E:114:THR:OG1	2.63	0.42
1:E:153:PHE:CD1	1:E:157:LEU:CD1	3.02	0.42
1:D:240:LEU:C	1:D:242:PRO:HD3	2.45	0.42
1:E:45:PHE:CE1	1:E:72:TYR:CE2	3.08	0.42
1:D:95:LEU:HD11	1:D:166:LEU:CD1	2.50	0.42
1:F:35:THR:HG21	1:F:241:PHE:CE2	2.55	0.42
1:F:122:MET:HB3	1:F:179:PHE:HB3	2.01	0.42
1:D:55:ASP:OD1	1:D:55:ASP:N	2.44	0.42
1:D:95:LEU:HD11	1:D:166:LEU:HD11	2.00	0.42
1:F:15:ALA:O	1:F:16:PRO:C	2.62	0.42
1:B:141:TYR:O	1:B:145:LYS:HG3	2.19	0.42
1:D:20:ILE:CG1	1:D:21:THR:N	2.83	0.42
1:D:50:GLU:OE1	1:D:50:GLU:C	2.62	0.42
1:F:137:ARG:NH2	6:F:301:G16:O3X	2.52	0.42
1:C:187:TYR:HA	1:C:190:ARG:NH1	2.34	0.42
1:F:69:LEU:HD23	1:F:187:TYR:CZ	2.55	0.42
1:E:187:TYR:O	1:E:190:ARG:HG2	2.18	0.42
1:D:217:PHE:CE2	1:D:224:GLY:C	2.98	0.42
1:E:99:CYS:O	1:E:103:ILE:HG12	2.19	0.42
1:C:8:ASP:O	1:C:12:THR:HB	2.20	0.41
1:E:102:TYR:CE1	1:E:153:PHE:CD2	3.08	0.41
1:E:143:LEU:HD23	1:E:143:LEU:HA	1.92	0.41
1:B:82:ASN:OD1	1:B:85:GLY:N	2.52	0.41
1:F:138:ILE:O	1:F:139:GLU:C	2.63	0.41
1:E:84:GLN:H	1:E:84:GLN:CD	2.28	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:40:VAL:HG23	1:F:64:PHE:CB	2.51	0.41
1:F:108:LEU:HG	1:F:109:PRO:HD2	2.03	0.41
1:F:137:ARG:HD3	1:F:173:GLN:OE1	2.20	0.41
1:F:154:VAL:HG21	1:F:170:ILE:HG13	2.02	0.41
1:F:232:ASP:OD1	1:F:235:ARG:NE	2.54	0.41
1:F:76:LYS:HD3	1:F:76:LYS:HA	1.81	0.41
1:C:232:ASP:OD1	1:C:235:ARG:NH2	2.49	0.41
1:B:125:VAL:O	1:B:176:ILE:N	2.53	0.41
1:D:27:PHE:CZ	1:D:238:GLU:HA	2.55	0.41
1:F:36:LYS:HD2	1:F:197:TYR:HE1	1.85	0.41
1:D:27:PHE:HE1	1:D:34:LYS:NZ	2.19	0.41
1:D:241:PHE:N	1:D:242:PRO:HD3	2.35	0.41
1:F:102:TYR:CE1	1:F:106:ILE:HD11	2.54	0.41
1:F:156:ASP:OD1	1:F:156:ASP:N	2.54	0.41
1:F:24:MET:HE1	1:F:230:PRO:O	2.21	0.41
1:C:83:ILE:HD12	1:C:180:PRO:HD3	2.02	0.41
1:B:56:VAL:HA	1:B:59:LYS:NZ	2.35	0.41
1:D:231:GLU:OE1	1:D:231:GLU:N	2.54	0.41
1:F:65:PRO:HD2	1:F:70:VAL:HG12	2.01	0.41
1:C:73:LYS:HB3	1:C:78:LEU:HD11	2.03	0.41
1:E:70:VAL:HG22	1:E:80:LYS:HG2	2.03	0.41
1:A:82:ASN:OD1	1:A:84:GLN:HG2	2.21	0.40
1:E:86:HIS:HB3	1:E:87:LEU:HD22	2.03	0.40
1:E:98:TYR:CD2	1:E:160:GLU:HG2	2.56	0.40
1:E:131:SER:O	1:E:132:CYS:C	2.63	0.40
1:D:106:ILE:HG22	1:D:108:LEU:HG	2.03	0.40
1:D:134:GLN:O	1:D:138:ILE:HD12	2.21	0.40
1:F:199:THR:HG21	1:F:201:TYR:CZ	2.55	0.40
1:F:40:VAL:HG23	1:F:64:PHE:HB3	2.02	0.40
1:E:124:ASN:HB3	1:E:130:ARG:HH21	1.86	0.40
1:C:3:THR:HG22	1:C:199:THR:HB	2.04	0.40
1:D:20:ILE:CG1	1:D:21:THR:H	2.35	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)
1:D:148:HIS:HE2	5:A:306:NA:NA[4_545]	1.53	0.07

5.3 Torsion angles

5.3.1 Protein backbone

There are no protein backbone outliers to report in this entry.

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	207/211 (98%)	204 (99%)	3 (1%)	59	75
1	B	210/211 (100%)	205 (98%)	5 (2%)	43	65
1	C	209/211 (99%)	204 (98%)	5 (2%)	43	65
1	D	204/211 (97%)	198 (97%)	6 (3%)	37	61
1	E	196/211 (93%)	193 (98%)	3 (2%)	57	74
1	F	189/211 (90%)	182 (96%)	7 (4%)	30	54
All	All	1215/1266 (96%)	1186 (98%)	29 (2%)	43	65

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	GLU
1	A	173	GLN
1	A	175	SER
1	B	21	THR
1	B	79	CYS
1	B	156	ASP
1	B	184	ASP
1	B	233	THR
1	D	14	THR
1	D	82	ASN
1	D	101	SER
1	D	149	ILE
1	D	175	SER
1	D	237	CYS
1	F	28	LEU

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Mol	Chain	Res	Type
1	F	78	LEU
1	F	79	CYS
1	F	101	SER
1	F	128	ILE
1	F	169	SER
1	F	227	VAL
1	C	79	CYS
1	C	82	ASN
1	C	93	GLN
1	C	131	SER
1	C	175	SER
1	E	119	ARG
1	E	130	ARG
1	E	131	SER

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

Mol	Chain	Res	Type
1	A	105	ASN
1	A	191	HIS
1	B	124	ASN
1	B	191	HIS
1	F	84	GLN
1	F	105	ASN
1	C	82	ASN
1	E	191	HIS

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

Of 22 ligands modelled in this entry, 17 are monoatomic - leaving 5 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
6	G16	F	301	-	19,20,20	1.57	2 (10%)	30,31,31	0.93	1 (3%)
6	G16	D	301	-	19,20,20	1.58	2 (10%)	30,31,31	0.99	1 (3%)
6	G16	C	301	-	19,20,20	1.73	4 (21%)	30,31,31	0.96	1 (3%)
2	GOL	A	301	-	5,5,5	0.70	0	5,5,5	1.00	0
6	G16	B	301	-	19,20,20	1.60	2 (10%)	30,31,31	1.09	2 (6%)

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
6	G16	F	301	-	-	0/11/31/31	0/1/1/1
6	G16	D	301	-	-	3/11/31/31	0/1/1/1
6	G16	C	301	-	-	3/11/31/31	0/1/1/1
2	GOL	A	301	-	-	0/4/4/4	-
6	G16	B	301	-	-	3/11/31/31	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	301	G16	P'-O1	4.82	1.68	1.59
6	F	301	G16	P'-O1	4.33	1.67	1.59
6	D	301	G16	P'-O1	4.28	1.67	1.59
6	B	301	G16	P'-O1	4.26	1.67	1.59
6	C	301	G16	O5-C1	4.11	1.52	1.41
6	D	301	G16	O5-C1	3.90	1.51	1.41
6	B	301	G16	O5-C1	3.75	1.51	1.41
6	F	301	G16	O5-C1	3.54	1.51	1.41
6	C	301	G16	O5-C5	2.22	1.49	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	301	G16	P-O6	2.02	1.66	1.60

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	301	G16	O5-C1-O1	-2.86	107.62	111.36
6	B	301	G16	O5-C5-C6	2.86	112.36	106.69
6	B	301	G16	C1-C2-C3	2.54	115.36	110.01
6	D	301	G16	O5-C5-C6	2.46	111.57	106.69
6	F	301	G16	C4-C3-C2	2.18	114.66	110.83

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	301	G16	C4-C5-C6-O6
6	B	301	G16	O5-C5-C6-O6
6	D	301	G16	C4-C5-C6-O6
6	D	301	G16	O5-C5-C6-O6
6	C	301	G16	C4-C5-C6-O6
6	C	301	G16	O5-C5-C6-O6
6	D	301	G16	C6-O6-P-O1P
6	C	301	G16	C1-O1-P'-O3X
6	B	301	G16	C2-C1-O1-P'

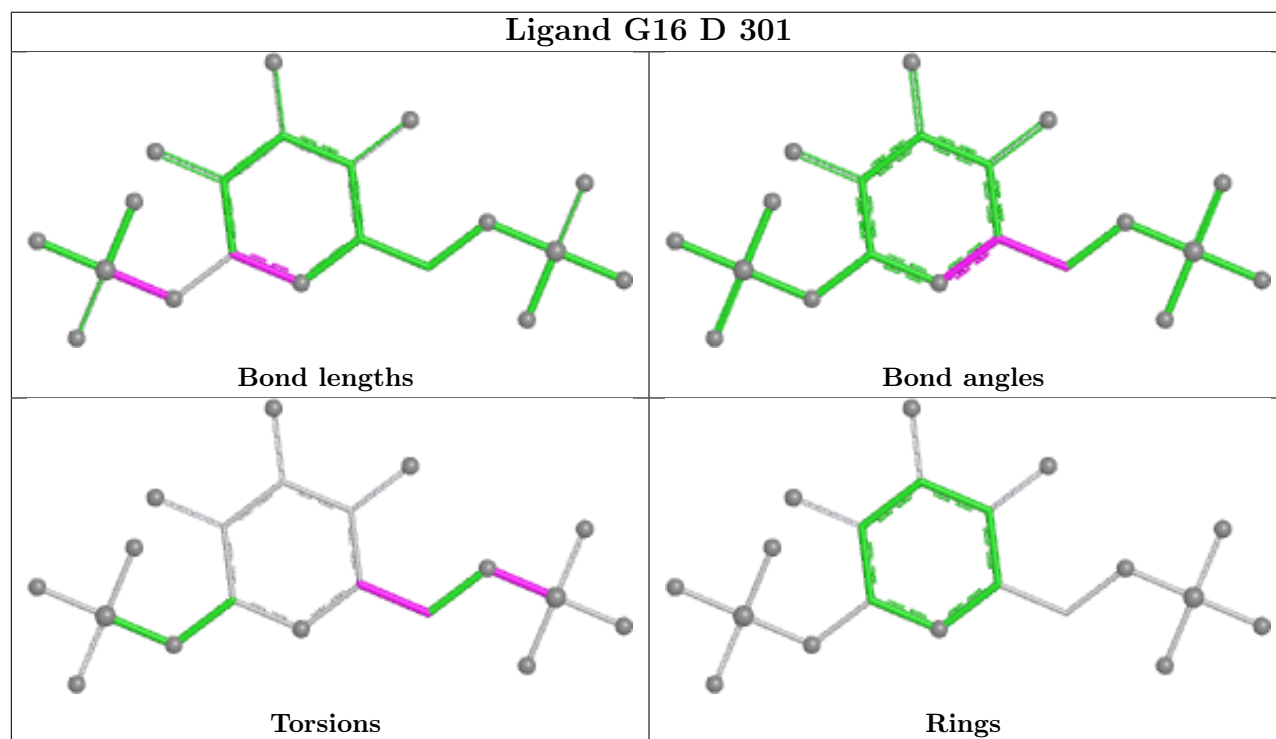
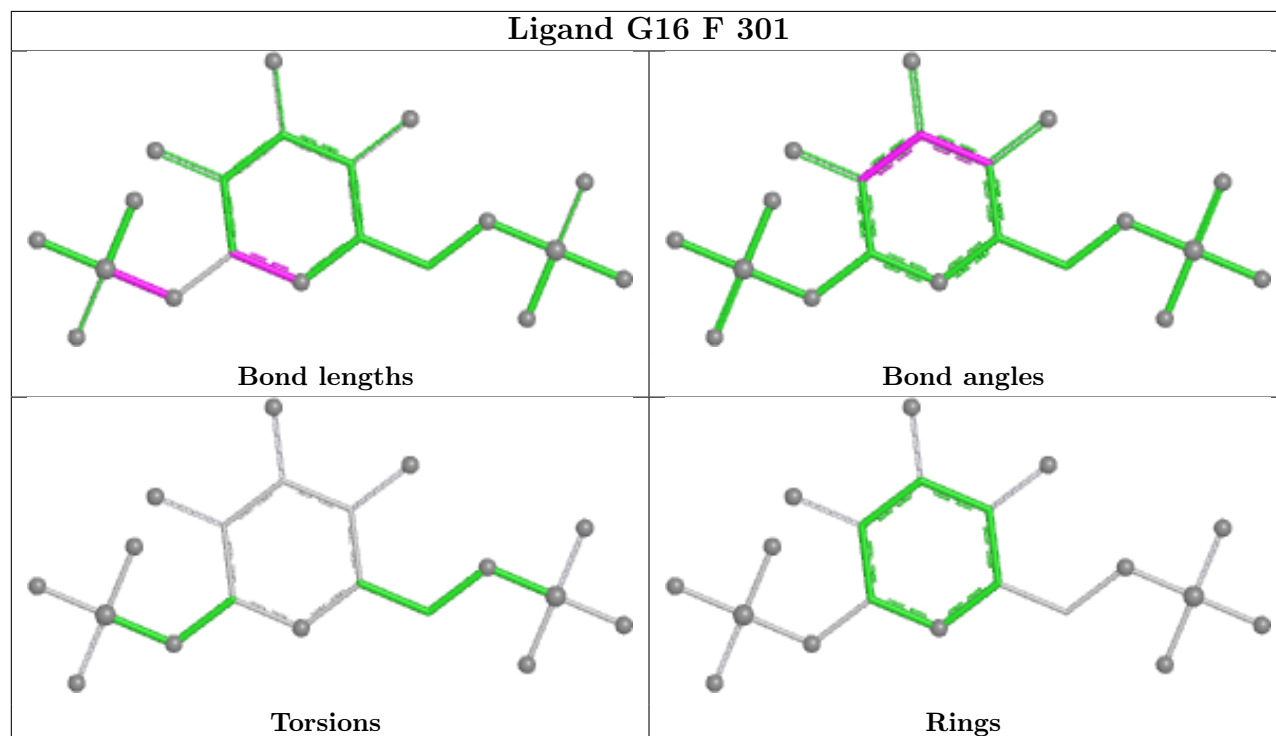
There are no ring outliers.

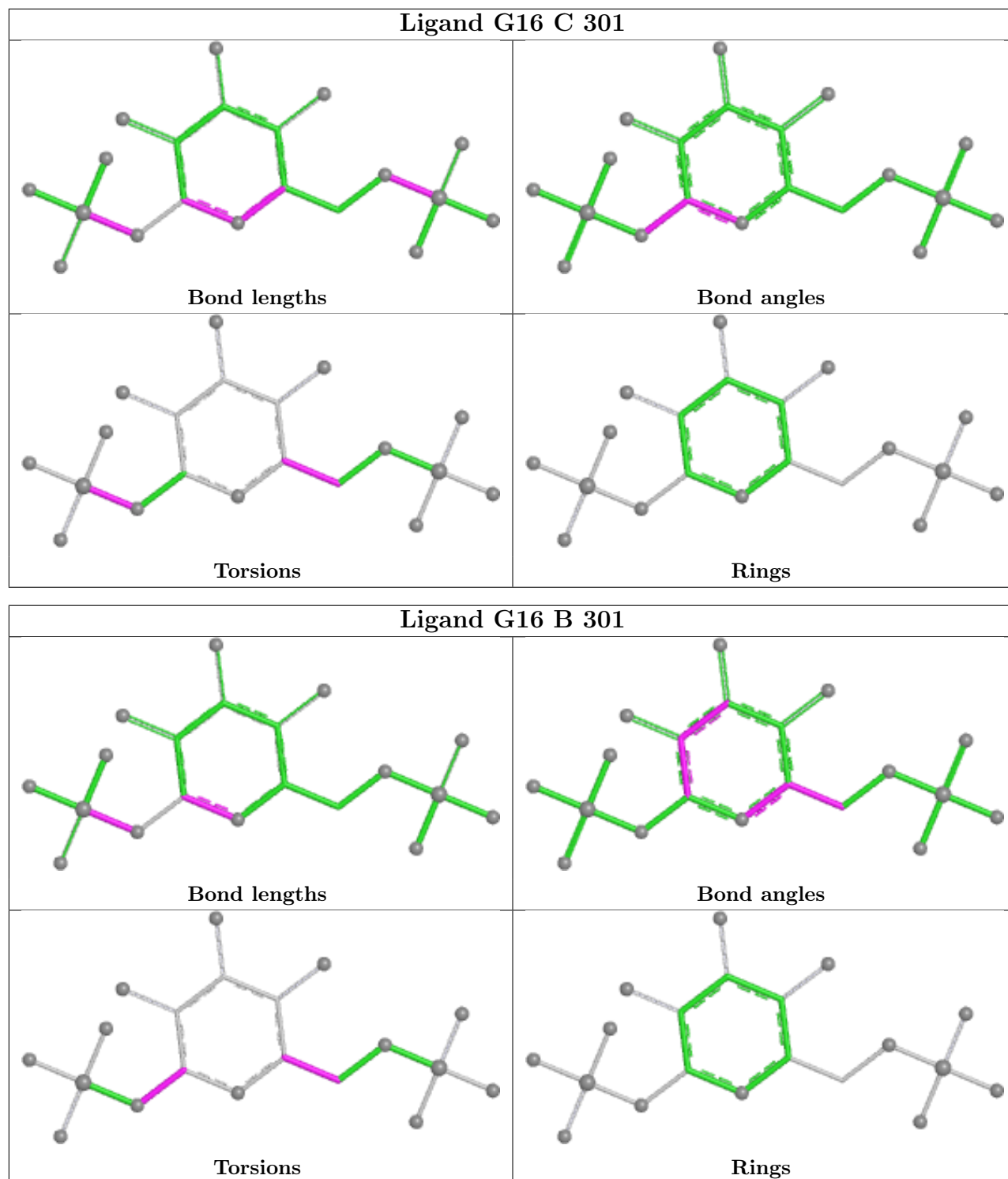
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	F	301	G16	1	0
6	C	301	G16	1	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 ⓘ

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	242/244 (99%)	-0.42	0 100 100	38, 54, 69, 87	0
1	B	243/244 (99%)	-0.13	1 (0%) 88 89	44, 64, 88, 103	0
1	C	243/244 (99%)	-0.18	0 100 100	46, 61, 83, 95	0
1	D	242/244 (99%)	0.03	2 (0%) 82 82	45, 69, 100, 117	0
1	E	242/244 (99%)	0.26	7 (2%) 53 52	52, 86, 110, 121	0
1	F	241/244 (98%)	0.40	11 (4%) 37 36	62, 90, 112, 133	0
All	All	1453/1464 (99%)	-0.01	21 (1%) 73 73	38, 68, 105, 133	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	1	MET	3.5
1	F	148	HIS	3.3
1	E	38	GLY	3.3
1	F	2	ALA	3.2
1	F	144	ASP	3.0
1	E	90	ASP	3.0
1	E	156	ASP	2.8
1	F	90	ASP	2.7
1	F	118	PHE	2.6
1	F	130	ARG	2.6
1	F	149	ILE	2.4
1	E	1	MET	2.3
1	E	157	LEU	2.3
1	E	108	LEU	2.2
1	F	143	LEU	2.1
1	F	110	LYS	2.1
1	B	194	HIS	2.1
1	D	30	LYS	2.1
1	E	127	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	F	123	LEU	2.1
1	F	147	GLU	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(Å ²)	Q<0.9
6	G16	C	301	20/20	0.61	0.13	75,100,119,120	30
6	G16	F	301	20/20	0.68	0.14	94,108,125,126	30
2	GOL	A	301	6/6	0.68	0.15	71,86,101,102	0
6	G16	B	301	20/20	0.74	0.12	63,85,99,108	30
3	MG	E	302	1/1	0.78	0.15	66,66,66,66	0
6	G16	D	301	20/20	0.79	0.10	57,80,96,97	30
4	CL	E	304	1/1	0.89	0.07	74,74,74,74	0
3	MG	A	304	1/1	0.91	0.11	56,56,56,56	0
4	CL	E	303	1/1	0.92	0.11	62,62,62,62	0
3	MG	F	303	1/1	0.93	0.06	65,65,65,65	0
3	MG	D	303	1/1	0.94	0.11	65,65,65,65	0
5	NA	A	306	1/1	0.96	0.08	57,57,57,57	0
4	CL	A	305	1/1	0.97	0.07	44,44,44,44	0
3	MG	C	303	1/1	0.97	0.04	35,35,35,35	0
3	MG	F	302	1/1	0.98	0.03	62,62,62,62	0
4	CL	B	303	1/1	0.98	0.08	60,60,60,60	0
3	MG	A	303	1/1	0.98	0.04	29,29,29,29	0
3	MG	A	302	1/1	0.99	0.04	48,48,48,48	0
3	MG	E	301	1/1	0.99	0.02	59,59,59,59	0
3	MG	B	302	1/1	0.99	0.02	52,52,52,52	0
3	MG	D	302	1/1	0.99	0.05	64,64,64,64	0

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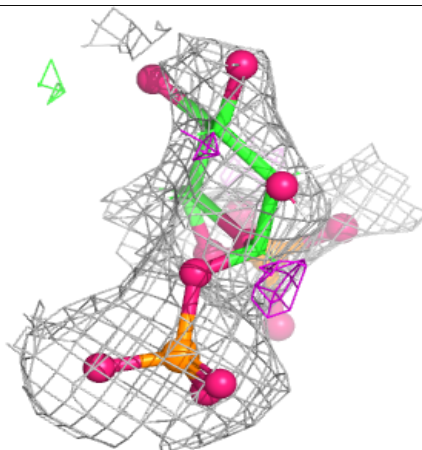
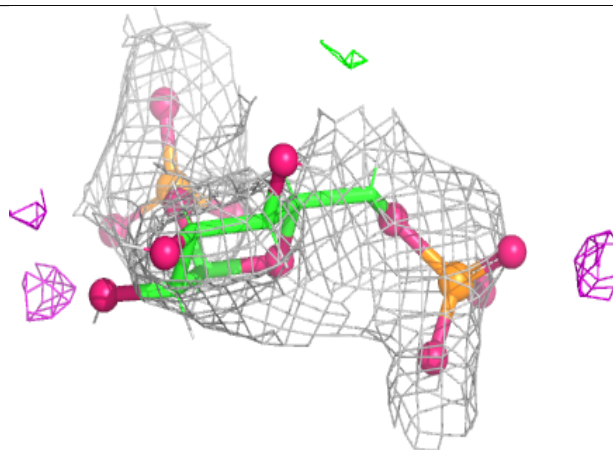
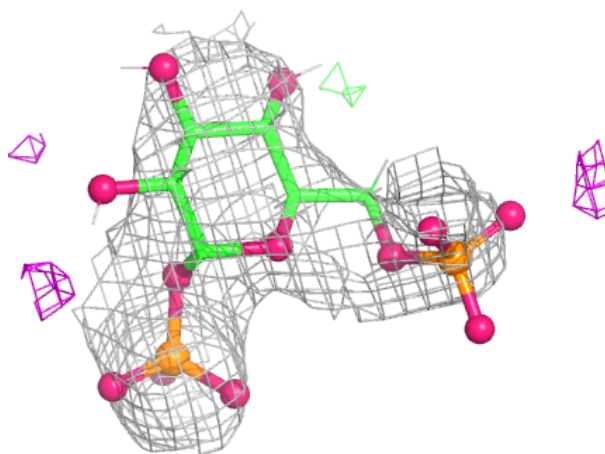
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	MG	C	302	1/1	1.00	0.03	49,49,49,49	0

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.

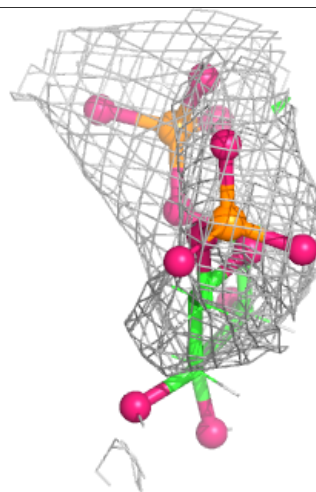
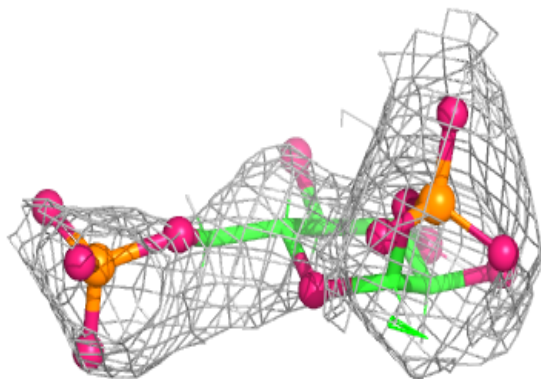
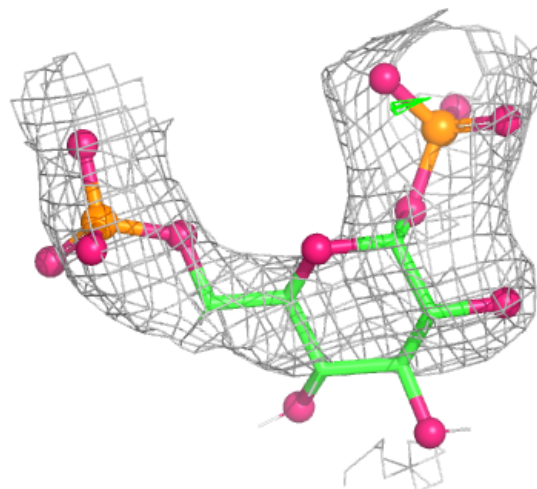
Electron density around G16 C 301:

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



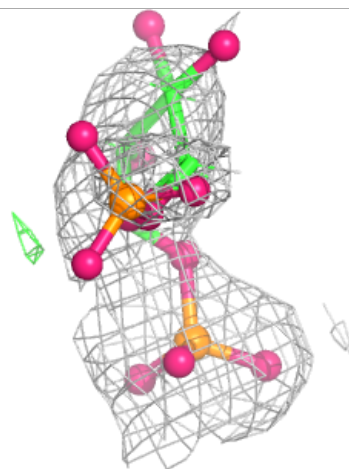
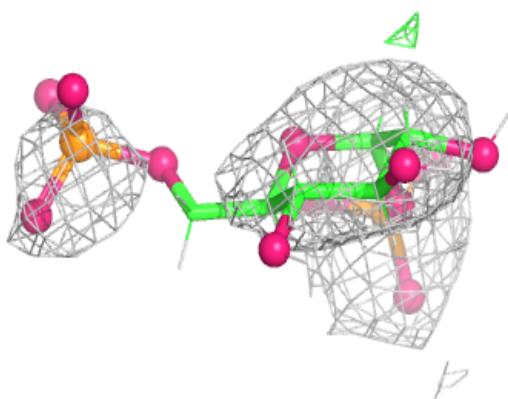
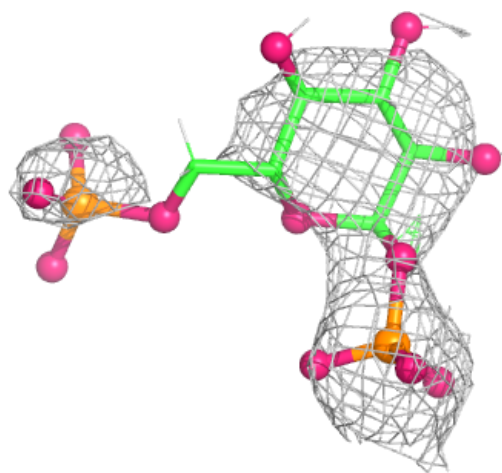
Electron density around G16 F 301:

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



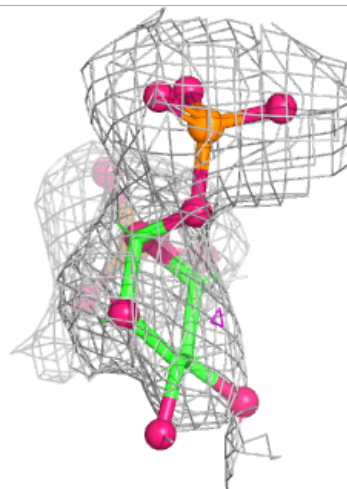
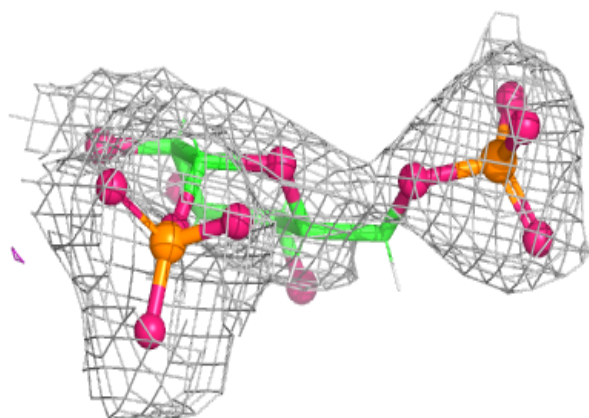
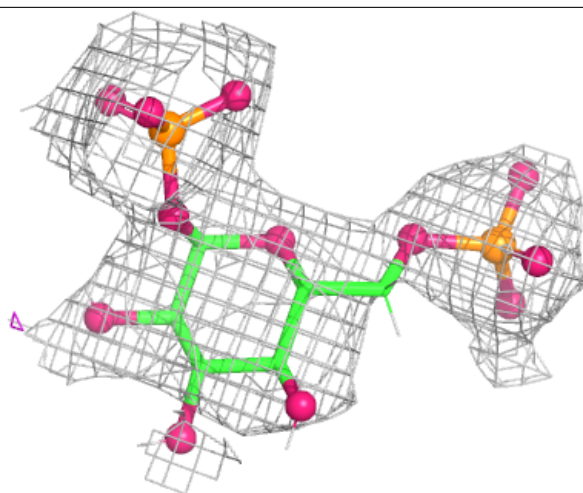
Electron density around G16 B 301:

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



Electron density around G16 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.