



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 06:40 AM UTC

PDB ID : 5U51 / pdb_00005u51
Title : Structure of Francisella tularensis heterodimeric SspA (MglA-SspA) in complex with ppGpp
Authors : Cuthbert, B.J.; Schumacher, M.A.; Brennan, R.G.
Deposited on : 2016-12-06
Resolution : 2.80 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

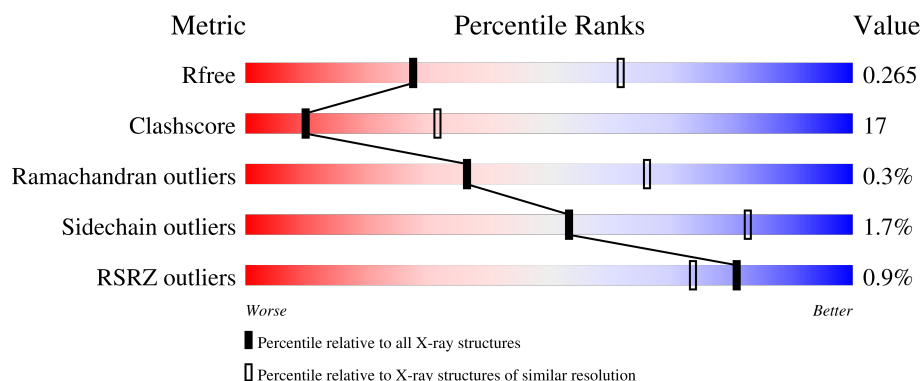
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

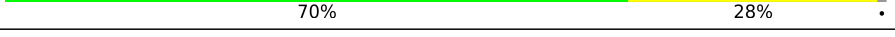
The reported resolution of this entry is 2.80 Å.

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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	C	211	
1	D	211	
2	A	204	
2	B	204	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6450 atoms, of which 87 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 Stringent starvation protein A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	195	Total	C	N	O	S	0	0	0
			1493	959	243	283	8			
1	D	194	Total	C	N	O	S	0	0	0
			1500	970	244	277	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	SER	-	expression tag	UNP A0A0E2ZL39
C	1	ASN	-	expression tag	UNP A0A0E2ZL39
C	2	ALA	-	expression tag	UNP A0A0E2ZL39
C	3	MET	-	expression tag	UNP A0A0E2ZL39
D	0	SER	-	expression tag	UNP A0A0E2ZL39
D	1	ASN	-	expression tag	UNP A0A0E2ZL39
D	2	ALA	-	expression tag	UNP A0A0E2ZL39
D	3	MET	-	expression tag	UNP A0A0E2ZL39

- Molecule 2 is a protein called Macrophage growth locus A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	203	Total	C	N	O	S	0	0	0
			1582	1031	253	293	5			
2	B	201	Total	C	N	O	S	0	0	0
			1609	1046	257	301	5			

There are 6 discrepancies between the modelled and reference sequences:

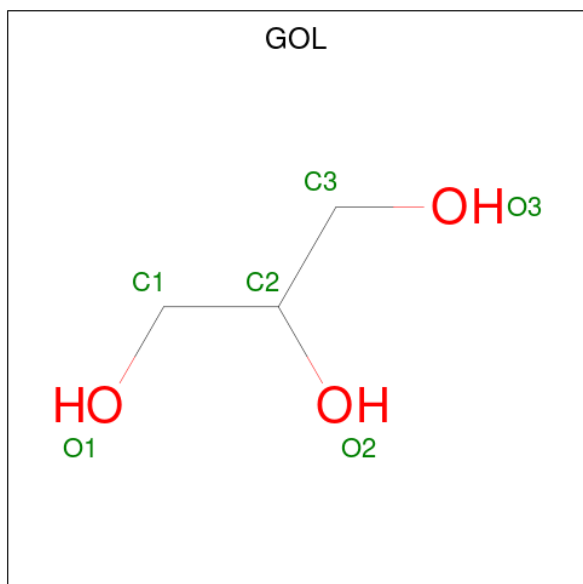
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP A0A0E2ZLH6
A	-1	ASN	-	expression tag	UNP A0A0E2ZLH6
A	0	ALA	-	expression tag	UNP A0A0E2ZLH6
B	-2	SER	-	expression tag	UNP A0A0E2ZLH6

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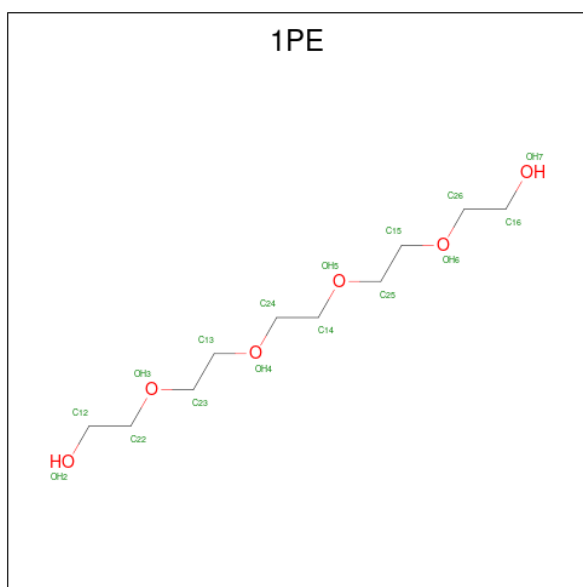
Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	ASN	-	expression tag	UNP A0A0E2ZLH6
B	0	ALA	-	expression tag	UNP A0A0E2ZLH6

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	C	1	Total	C	O		0	0
			6	3	3			
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 4 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).

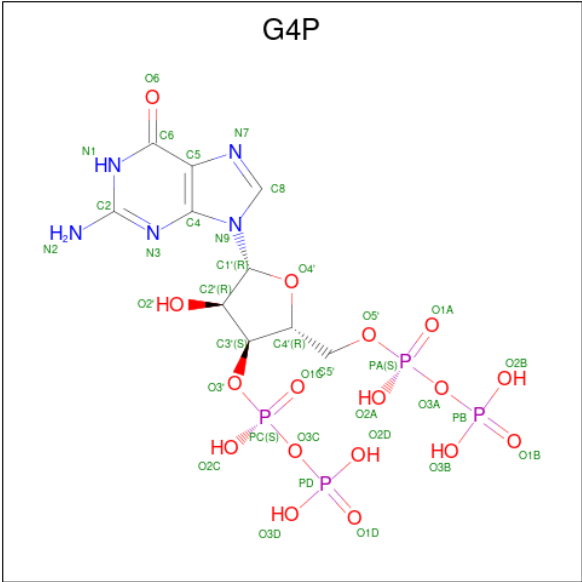


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	H	O	0	0
			38	10	22	6		
4	B	1	Total	C	H	O	0	0
			38	10	22	6		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	1	Total	Mg	0	0
			1	1		
5	B	1	Total	Mg	0	0
			1	1		

- Molecule 6 is GUANOSINE-5',3'-TETRAPHOSPHATE (CCD ID: G4P) (formula: C₁₀H₁₇N₅O₁₇P₄).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			36	10	5	17	4		
6	B	1	Total	C	H	N	O	0	0
			47	10	11	5	17		

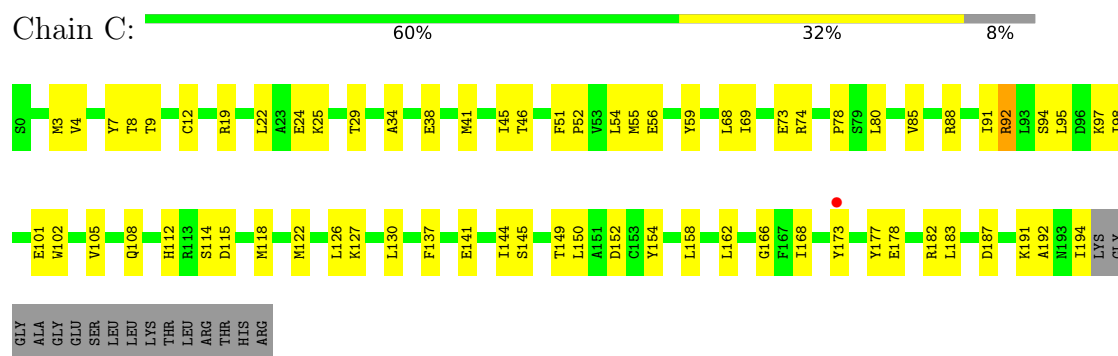
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	11	Total	O	0	0
			11	11		
7	A	8	Total	O	0	0
			8	8		
7	D	11	Total	O	0	0
			11	11		
7	B	13	Total	O	0	0
			13	13		

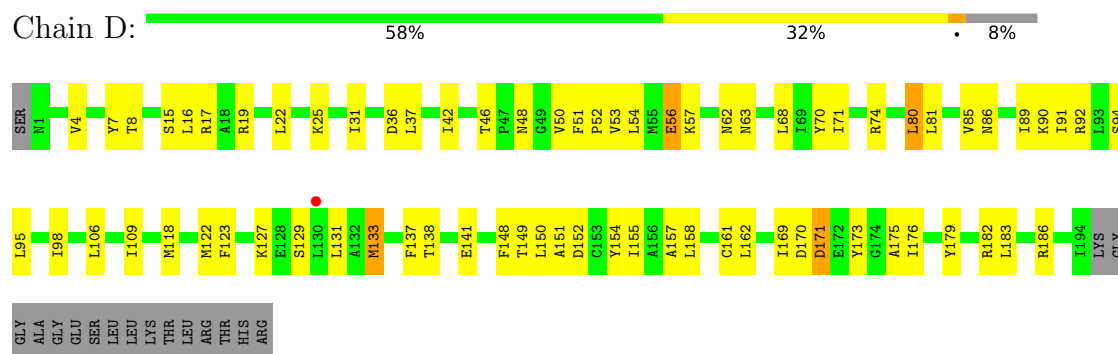
3 Residue-property plots [i](#)

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.

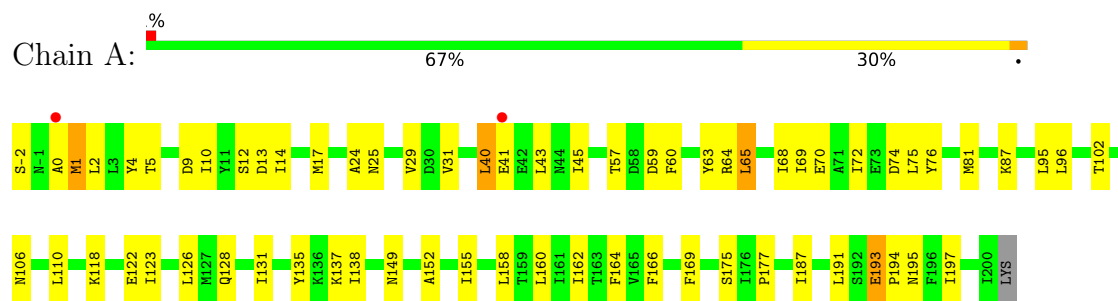
- Molecule 1: Stringent starvation protein A



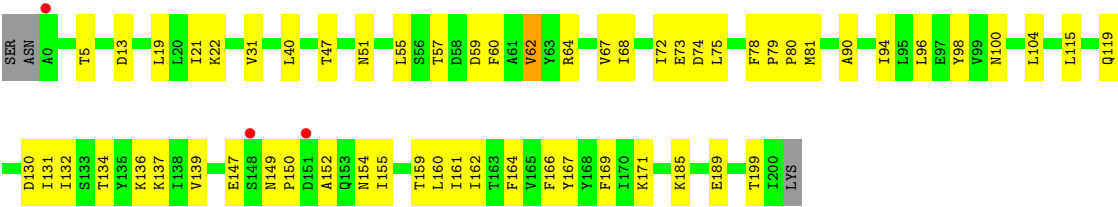
- Molecule 1: Stringent starvation protein A



- Molecule 2: Macrophage growth locus A



- Molecule 2: Macrophage growth locus A



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.62Å 111.56Å 141.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	87.61 – 2.80 87.61 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (87.61-2.80) 99.9 (87.61-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 2.82Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.212 , 0.265 0.212 , 0.265	Depositor DCC
R_{free} test set	1188 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	58.7	Xtriage
Anisotropy	0.440	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 37.5	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	0.93	EDS
Total number of atoms	6450	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, G4P, 1PE, GOL

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	C	0.15	0/1521	0.35	0/2066
1	D	0.14	0/1529	0.32	0/2075
2	A	0.14	0/1610	0.33	0/2194
2	B	0.13	0/1637	0.32	0/2223
All	All	0.14	0/6297	0.33	0/8558

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1493	0	1445	55	1
1	D	1500	0	1475	57	0
2	A	1582	0	1598	64	1
2	B	1609	0	1656	57	0
3	A	6	8	8	0	0
3	C	24	24	32	0	0
4	B	16	22	22	1	0
4	C	16	22	22	0	0
5	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	1	0	0	0	0
6	A	36	0	10	1	0
6	B	36	11	10	0	0
7	A	8	0	0	1	0
7	B	13	0	0	0	0
7	C	11	0	0	0	0
7	D	11	0	0	0	0
All	All	6363	87	6278	220	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 17.

All (220) 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)
2:A:57:THR:HG23	2:A:59:ASP:H	1.32	0.92
2:A:193:GLU:HG2	2:A:194:PRO:HD2	1.52	0.91
2:A:31:VAL:HG13	2:A:40:LEU:HB2	1.56	0.88
2:A:64:ARG:NH1	6:A:301:G4P:O2A	2.08	0.86
1:D:56:GLU:OE1	1:D:70:TYR:OH	1.94	0.85
1:D:19:ARG:HD3	1:D:31:ILE:HD13	1.60	0.83
1:C:4:VAL:HG12	1:C:56:GLU:HG2	1.60	0.83
1:D:25:LYS:HE2	1:D:80:LEU:HD22	1.60	0.83
2:A:138:ILE:HD12	2:A:162:ILE:HD12	1.59	0.82
2:B:47:THR:HG21	2:B:51:ASN:H	1.45	0.81
2:A:81:MET:HE2	2:A:81:MET:HA	1.61	0.80
2:B:81:MET:HA	2:B:81:MET:HE2	1.67	0.77
2:A:95:LEU:HD23	2:A:158:LEU:HD13	1.68	0.76
1:C:4:VAL:CG2	1:C:29:THR:HG22	2.17	0.75
1:D:25:LYS:HE2	1:D:80:LEU:CD2	2.18	0.74
2:A:193:GLU:O	2:A:197:ILE:HD12	1.90	0.72
1:C:3:MET:HE2	1:C:74:ARG:HH22	1.54	0.72
2:A:-2:SER:N	7:A:401:HOH:O	2.22	0.72
2:A:17:MET:HG2	2:A:160:LEU:HD11	1.70	0.71
1:C:115:ASP:HB3	1:C:118:MET:CB	2.20	0.71
1:D:118:MET:HG3	1:D:122:MET:HE2	1.71	0.71
1:D:63:ASN:HD22	2:B:94:ILE:HG12	1.55	0.70
1:D:19:ARG:CD	1:D:31:ILE:HD13	2.22	0.70
2:A:187:ILE:O	2:A:191:LEU:HD12	1.91	0.69
1:D:137:PHE:HD2	1:D:176:ILE:HG12	1.57	0.69
2:B:57:THR:HG23	2:B:59:ASP:H	1.57	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:53:VAL:HG22	1:D:62:ASN:HB3	1.73	0.69
2:A:95:LEU:HB3	2:A:158:LEU:HD11	1.75	0.69
1:D:85:VAL:HG11	2:B:75:LEU:HD23	1.75	0.69
2:A:149:ASN:ND2	2:A:152:ALA:HB2	2.07	0.68
1:C:102:TRP:O	1:C:105:VAL:HG12	1.94	0.68
1:C:158:LEU:O	1:C:162:LEU:HD23	1.94	0.67
2:B:149:ASN:ND2	2:B:152:ALA:HB2	2.10	0.67
1:D:179:TYR:CE2	1:D:183:LEU:HD11	2.30	0.66
2:A:31:VAL:CG1	2:A:40:LEU:HD22	2.26	0.66
1:D:179:TYR:CZ	1:D:183:LEU:HD11	2.31	0.66
1:D:7:TYR:CD1	1:D:37:LEU:HD21	2.31	0.66
2:B:73:GLU:OE2	2:B:80:PRO:HA	1.97	0.65
2:B:31:VAL:HG21	2:B:40:LEU:HD13	1.77	0.64
2:A:95:LEU:HD23	2:A:158:LEU:CD1	2.28	0.64
1:C:166:GLY:O	1:C:168:ILE:HD12	1.98	0.64
2:A:81:MET:CE	2:A:155:ILE:HB	2.28	0.64
2:B:19:LEU:HD21	2:B:72:ILE:HD13	1.79	0.64
2:A:40:LEU:HD12	2:A:40:LEU:O	1.98	0.64
2:A:81:MET:HE1	2:A:155:ILE:HD12	1.80	0.64
1:C:127:LYS:HG3	1:C:173:TYR:OH	1.99	0.63
2:A:5:THR:HB	2:A:12:SER:HB3	1.81	0.63
1:D:138:THR:HG22	1:D:175:ALA:CB	2.29	0.63
1:D:106:LEU:HD13	1:D:162:LEU:HD23	1.80	0.62
1:D:109:ILE:HD13	1:D:123:PHE:HE1	1.64	0.62
2:A:81:MET:HE2	2:A:155:ILE:HB	1.82	0.62
1:D:170:ASP:H	1:D:173:TYR:HE2	1.46	0.61
1:C:3:MET:HE2	1:C:74:ARG:NH2	2.16	0.61
2:B:5:THR:O	2:B:31:VAL:HG12	2.01	0.61
1:C:118:MET:O	1:C:122:MET:HG3	2.00	0.60
2:A:40:LEU:HD12	2:A:40:LEU:C	2.25	0.60
1:C:187:ASP:OD1	1:C:191:LYS:NZ	2.35	0.60
2:B:131:ILE:HD11	2:B:166:PHE:CE1	2.36	0.60
2:A:131:ILE:HD11	2:A:166:PHE:CE1	2.36	0.60
1:D:86:ASN:O	1:D:90:LYS:HD3	2.02	0.60
2:B:167:TYR:CZ	2:B:171:LYS:HD2	2.37	0.60
1:D:92:ARG:NH1	2:B:74:ASP:OD2	2.35	0.59
2:A:4:TYR:CG	2:A:43:LEU:HD13	2.37	0.59
1:C:141:GLU:O	1:C:182:ARG:NH1	2.21	0.59
1:C:126:LEU:O	1:C:130:LEU:HD13	2.02	0.58
2:A:118:LYS:O	2:A:122:GLU:HG3	2.03	0.58
1:C:173:TYR:O	1:C:177:TYR:HB2	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:173:TYR:O	1:C:177:TYR:CB	2.52	0.57
2:B:130:ASP:O	2:B:134:THR:HG23	2.05	0.57
2:B:55:LEU:HB3	2:B:62:VAL:HG13	1.87	0.57
1:D:138:THR:HG22	1:D:175:ALA:HB2	1.86	0.57
1:C:137:PHE:HZ	1:C:144:ILE:H	1.52	0.56
2:A:2:LEU:HD11	2:A:29:VAL:HG21	1.86	0.56
1:C:4:VAL:HG22	1:C:29:THR:HG22	1.87	0.56
1:C:144:ILE:O	1:C:145:SER:OG	2.22	0.56
2:B:19:LEU:CD2	2:B:72:ILE:HD13	2.36	0.56
1:C:92:ARG:HH21	2:A:70:GLU:HB3	1.70	0.56
1:C:178:GLU:O	1:C:182:ARG:HG3	2.06	0.55
2:A:135:TYR:HD1	2:A:162:ILE:HD13	1.71	0.55
1:C:92:ARG:HE	2:A:70:GLU:HB3	1.71	0.55
2:B:185:LYS:O	2:B:189:GLU:HG2	2.05	0.55
1:D:85:VAL:CG1	2:B:75:LEU:HD23	2.37	0.55
1:C:149:THR:HG22	1:C:150:LEU:N	2.20	0.55
2:B:47:THR:CG2	2:B:51:ASN:H	2.17	0.55
1:C:85:VAL:HG13	2:A:74:ASP:HB3	1.89	0.54
1:D:137:PHE:CD2	1:D:176:ILE:HG12	2.40	0.54
2:B:31:VAL:CG2	2:B:40:LEU:HD13	2.36	0.54
1:C:94:SER:O	1:C:98:ILE:HG13	2.07	0.54
1:C:88:ARG:HG2	1:C:92:ARG:NH1	2.22	0.54
1:C:69:ILE:O	1:C:73:GLU:HG3	2.08	0.54
1:C:88:ARG:HG2	1:C:92:ARG:HH12	1.73	0.54
2:A:13:ASP:HB3	2:A:164:PHE:CE1	2.43	0.53
2:A:149:ASN:CG	2:A:152:ALA:HB2	2.34	0.53
1:D:46:THR:HB	1:D:53:VAL:HG21	1.91	0.53
2:A:102:THR:O	2:A:106:ASN:ND2	2.37	0.53
2:B:90:ALA:O	2:B:94:ILE:HG13	2.07	0.53
1:C:149:THR:HG22	1:C:150:LEU:H	1.74	0.53
1:D:51:PHE:HB3	1:D:52:PRO:HA	1.90	0.53
1:C:24:GLU:HG2	1:C:80:LEU:HD21	1.91	0.52
2:B:62:VAL:HG22	2:B:68:ILE:HG12	1.90	0.52
1:D:4:VAL:HG12	1:D:56:GLU:HG3	1.90	0.52
2:B:47:THR:HG21	2:B:51:ASN:N	2.20	0.52
1:C:45:ILE:HG23	1:C:55:MET:HE3	1.91	0.52
1:D:158:LEU:O	1:D:162:LEU:HG	2.09	0.52
1:D:37:LEU:HB2	1:D:42:ILE:HD11	1.92	0.52
1:C:97:LYS:HE3	2:A:63:TYR:OH	2.10	0.52
1:C:97:LYS:HD2	1:C:101:GLU:OE1	2.10	0.52
1:D:46:THR:HG23	1:D:46:THR:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:128:GLN:OE1	2:A:175:SER:HB3	2.10	0.52
1:D:80:LEU:HB3	1:D:150:LEU:HG	1.92	0.52
2:A:1:MET:HB2	2:A:25:ASN:O	2.10	0.51
2:A:131:ILE:HD12	2:A:169:PHE:CZ	2.45	0.51
1:D:95:LEU:HD22	1:D:154:TYR:CE2	2.45	0.51
2:B:80:PRO:HG2	2:B:154:ASN:CB	2.39	0.51
1:D:129:SER:O	1:D:133:MET:HG3	2.10	0.51
2:A:31:VAL:CG1	2:A:40:LEU:HB2	2.37	0.51
1:C:152:ASP:HB3	1:C:183:LEU:HD11	1.93	0.50
2:A:4:TYR:CD1	2:A:43:LEU:HD13	2.47	0.50
1:D:15:SER:HA	1:D:52:PRO:HG3	1.92	0.50
1:D:54:LEU:HD23	1:D:54:LEU:C	2.35	0.50
2:A:96:LEU:O	2:A:96:LEU:HD23	2.12	0.50
1:D:8:THR:HB	1:D:15:SER:HB3	1.94	0.50
1:D:94:SER:O	1:D:98:ILE:HG13	2.11	0.50
2:A:193:GLU:OE2	2:A:195:ASN:HB2	2.11	0.50
2:B:131:ILE:HD12	2:B:169:PHE:CZ	2.47	0.50
1:D:48:ASN:HB2	1:D:50:VAL:HG23	1.93	0.49
1:D:56:GLU:OE2	1:D:74:ARG:NE	2.43	0.49
2:A:175:SER:O	2:A:177:PRO:HD3	2.12	0.49
1:C:51:PHE:HB3	1:C:52:PRO:HA	1.94	0.49
1:C:68:LEU:HD22	1:C:154:TYR:HE2	1.78	0.49
1:D:127:LYS:O	1:D:131:LEU:HG	2.13	0.49
2:B:64:ARG:O	2:B:68:ILE:HG13	2.12	0.49
2:B:68:ILE:O	2:B:72:ILE:HG13	2.12	0.49
1:C:19:ARG:NH1	1:C:192:ALA:HA	2.27	0.49
1:D:25:LYS:CE	1:D:80:LEU:HD22	2.38	0.49
1:C:4:VAL:HG23	1:C:29:THR:HG22	1.94	0.49
2:B:19:LEU:HD21	2:B:72:ILE:CD1	2.43	0.49
2:B:131:ILE:HD11	2:B:166:PHE:CD1	2.47	0.48
1:C:9:THR:HG23	1:C:12:CYS:HB2	1.95	0.48
1:D:171:ASP:OD1	1:D:171:ASP:N	2.46	0.48
2:A:64:ARG:O	2:A:68:ILE:HG13	2.14	0.48
2:B:100:ASN:HD22	2:B:104:LEU:HD12	1.79	0.48
2:A:68:ILE:O	2:A:72:ILE:HG13	2.14	0.48
2:B:62:VAL:HG23	2:B:67:VAL:HG12	1.97	0.47
1:D:169:ILE:HD12	1:D:169:ILE:N	2.30	0.47
2:A:41:GLU:O	2:A:45:ILE:HG13	2.14	0.47
2:B:13:ASP:HB3	2:B:164:PHE:CE1	2.50	0.47
2:B:57:THR:HG21	2:B:60:PHE:CE2	2.50	0.47
1:C:54:LEU:C	1:C:54:LEU:HD23	2.40	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:2:LEU:HD11	2:A:29:VAL:CG2	2.45	0.46
2:A:81:MET:HE1	2:A:155:ILE:HB	1.96	0.46
1:C:173:TYR:O	1:C:177:TYR:HB3	2.15	0.46
2:A:160:LEU:C	2:A:160:LEU:HD23	2.41	0.46
2:B:78:PHE:HA	2:B:79:PRO:HA	1.76	0.46
2:B:22:LYS:NZ	2:B:81:MET:HG2	2.31	0.46
2:A:10:ILE:O	2:A:14:ILE:HG13	2.16	0.46
2:B:81:MET:CE	2:B:155:ILE:HB	2.47	0.46
1:C:91:ILE:O	1:C:95:LEU:HG	2.15	0.45
1:D:149:THR:HG23	1:D:151:ALA:H	1.81	0.45
1:C:59:TYR:HB2	2:A:87:LYS:HD2	1.98	0.45
2:A:118:LYS:HE3	2:A:118:LYS:HB2	1.75	0.45
2:A:81:MET:HE1	2:A:155:ILE:CD1	2.47	0.45
1:C:46:THR:HG23	1:C:46:THR:O	2.17	0.45
1:D:16:LEU:HB2	1:D:157:ALA:HB1	1.99	0.45
2:A:57:THR:HG21	2:A:60:PHE:CE2	2.52	0.44
1:C:7:TYR:CE1	1:C:45:ILE:HD13	2.53	0.44
2:B:47:THR:HG21	2:B:51:ASN:HB2	1.98	0.44
1:D:109:ILE:HD13	1:D:123:PHE:CE1	2.50	0.44
1:D:141:GLU:O	1:D:182:ARG:NH1	2.50	0.44
1:C:68:LEU:HD22	1:C:154:TYR:CE2	2.53	0.44
2:B:96:LEU:HD12	2:B:161:ILE:HD12	2.00	0.44
2:B:147:GLU:O	2:B:150:PRO:HD3	2.18	0.44
2:A:131:ILE:HD11	2:A:166:PHE:CD1	2.52	0.44
1:D:106:LEU:HD21	1:D:161:CYS:HB3	1.99	0.44
1:D:148:PHE:CE1	1:D:186:ARG:HD3	2.53	0.44
1:D:149:THR:HG23	1:D:151:ALA:N	2.33	0.43
1:C:97:LYS:CE	2:A:63:TYR:OH	2.65	0.43
2:B:21:ILE:HG21	2:B:155:ILE:CD1	2.48	0.43
2:B:115:LEU:HD11	2:B:119:GLN:HG2	2.01	0.43
1:C:85:VAL:HG11	2:A:75:LEU:HD23	1.99	0.43
2:B:137:LYS:HD3	4:B:302:1PE:H132	2.00	0.43
1:C:194:ILE:HD12	1:C:194:ILE:C	2.44	0.43
2:A:9:ASP:OD2	2:A:12:SER:OG	2.30	0.43
2:A:40:LEU:C	2:A:40:LEU:CD1	2.91	0.43
1:D:151:ALA:O	1:D:155:ILE:HG13	2.19	0.43
2:B:199:THR:HG22	2:B:199:THR:O	2.19	0.42
1:D:22:LEU:HG	1:D:71:ILE:HD13	2.01	0.42
2:B:115:LEU:HD11	2:B:119:GLN:CG	2.49	0.42
2:A:24:ALA:HB1	2:A:76:TYR:CZ	2.54	0.42
1:D:36:ASP:C	1:D:37:LEU:HD12	2.45	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:63:ASN:ND2	2:B:94:ILE:HG12	2.29	0.42
2:B:167:TYR:CE2	2:B:171:LYS:HD2	2.54	0.42
1:C:19:ARG:HH12	1:C:192:ALA:HA	1.84	0.42
1:D:91:ILE:O	1:D:95:LEU:HG	2.19	0.42
2:B:31:VAL:HG22	2:B:31:VAL:O	2.20	0.42
2:A:57:THR:CG2	2:A:60:PHE:H	2.33	0.42
2:B:62:VAL:CG2	2:B:67:VAL:HG12	2.50	0.42
2:B:81:MET:HE1	2:B:155:ILE:HB	2.02	0.42
2:B:98:TYR:HE2	2:B:134:THR:CG2	2.33	0.42
1:C:22:LEU:HD13	1:C:29:THR:HG21	2.02	0.42
1:D:95:LEU:HD22	1:D:154:TYR:CZ	2.55	0.41
2:A:110:LEU:HD21	2:A:123:ILE:HB	2.02	0.41
2:A:57:THR:HG23	2:A:59:ASP:N	2.16	0.41
1:D:152:ASP:HB3	1:D:183:LEU:HD21	2.02	0.41
2:B:98:TYR:HE2	2:B:134:THR:HG21	1.85	0.41
1:C:25:LYS:HA	1:C:78:PRO:HD2	2.03	0.41
2:A:0:ALA:O	2:A:1:MET:C	2.63	0.41
2:A:122:GLU:O	2:A:126:LEU:HG	2.20	0.41
1:C:112:HIS:C	1:C:114:SER:H	2.28	0.41
1:C:41:MET:HE2	1:C:45:ILE:HD11	2.03	0.41
2:B:81:MET:HE1	2:B:155:ILE:HD12	2.03	0.41
1:C:8:THR:O	1:C:34:ALA:N	2.49	0.41
1:D:17:ARG:HG2	1:D:68:LEU:HD21	2.03	0.41
1:D:89:ILE:CD1	2:B:75:LEU:HD21	2.51	0.41
2:B:136:LYS:HA	2:B:139:VAL:HG12	2.03	0.41
1:C:108:GLN:HB3	1:C:122:MET:HE1	2.02	0.40
2:A:65:LEU:HD22	2:A:69:ILE:HG13	2.03	0.40
2:B:159:THR:HA	2:B:162:ILE:HG22	2.03	0.40
2:B:160:LEU:C	2:B:160:LEU:HD23	2.46	0.40
2:B:132:ILE:HD13	2:B:132:ILE:HA	1.96	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:C:114:SER:OG	2:A:137:LYS:NZ[4_555]	2.16	0.04

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	193/211 (92%)	185 (96%)	8 (4%)	0	100	100
1	D	192/211 (91%)	183 (95%)	8 (4%)	1 (0%)	24	55
2	A	201/204 (98%)	196 (98%)	4 (2%)	1 (0%)	24	55
2	B	199/204 (98%)	194 (98%)	5 (2%)	0	100	100
All	All	785/830 (95%)	758 (97%)	25 (3%)	2 (0%)	36	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	1	MET
1	D	57	LYS

5.3.2 Protein sidechains [i](#)

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	C	153/184 (83%)	151 (99%)	2 (1%)	61	86
1	D	154/184 (84%)	149 (97%)	5 (3%)	34	70
2	A	173/192 (90%)	170 (98%)	3 (2%)	53	83
2	B	182/192 (95%)	181 (100%)	1 (0%)	81	93
All	All	662/752 (88%)	651 (98%)	11 (2%)	53	83

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	38	GLU
1	C	92	ARG
2	A	40	LEU
2	A	65	LEU
2	A	193	GLU
1	D	56	GLU
1	D	80	LEU
1	D	81	LEU
1	D	133	MET
1	D	171	ASP
2	B	62	VAL

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

Mol	Chain	Res	Type
1	C	62	ASN
2	A	105	GLN
2	A	154	ASN
1	D	63	ASN
1	D	86	ASN
2	B	100	ASN
2	B	149	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 ⓘ

Of 11 ligands modelled in this entry, 2 are monoatomic - leaving 9 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
3	GOL	C	302	-	5,5,5	0.40	0	5,5,5	0.20	0
3	GOL	C	303	-	5,5,5	0.38	0	5,5,5	0.35	0
6	G4P	B	301	5	36,38,38	3.27	16 (44%)	56,61,61	1.90	13 (23%)
6	G4P	A	301	5	36,38,38	3.28	15 (41%)	56,61,61	1.88	10 (17%)
3	GOL	C	301	-	5,5,5	0.45	0	5,5,5	0.16	0
3	GOL	A	302	-	5,5,5	0.42	0	5,5,5	0.12	0
3	GOL	C	304	-	5,5,5	0.38	0	5,5,5	0.20	0
4	1PE	B	302	-	15,15,15	0.54	0	14,14,14	0.23	0
4	1PE	C	305	-	15,15,15	0.53	0	14,14,14	0.27	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	C	302	-	-	3/4/4/4	-
3	GOL	C	303	-	-	4/4/4/4	-
6	G4P	B	301	5	-	7/27/43/43	0/3/3/3
6	G4P	A	301	5	-	8/27/43/43	0/3/3/3
3	GOL	C	301	-	-	4/4/4/4	-
3	GOL	A	302	-	-	4/4/4/4	-
3	GOL	C	304	-	-	1/4/4/4	-
4	1PE	B	302	-	-	5/13/13/13	-
4	1PE	C	305	-	-	8/13/13/13	-

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	301	G4P	C3'-C4'	-8.52	1.30	1.52
6	B	301	G4P	C3'-C4'	-8.23	1.31	1.52
6	A	301	G4P	O4'-C4'	8.02	1.62	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	301	G4P	O4'-C4'	7.60	1.61	1.45
6	B	301	G4P	C4-N3	6.70	1.49	1.34
6	A	301	G4P	C4-N3	6.61	1.49	1.34
6	A	301	G4P	PA-O3A	6.03	1.66	1.59
6	B	301	G4P	PA-O3A	6.02	1.66	1.59
6	A	301	G4P	C2-N3	5.61	1.46	1.33
6	B	301	G4P	PC-O3C	5.55	1.65	1.59
6	B	301	G4P	C2-N3	5.52	1.46	1.33
6	A	301	G4P	PC-O3C	5.30	1.65	1.59
6	B	301	G4P	C2-N2	4.98	1.45	1.34
6	A	301	G4P	C2-N2	4.78	1.45	1.34
6	B	301	G4P	O4'-C1'	-4.74	1.31	1.42
6	A	301	G4P	O4'-C1'	-4.64	1.31	1.42
6	B	301	G4P	C2-N1	2.98	1.44	1.37
6	B	301	G4P	C5-N7	-2.86	1.33	1.39
6	A	301	G4P	C5-N7	-2.75	1.33	1.39
6	A	301	G4P	C2-N1	2.68	1.44	1.37
6	A	301	G4P	O2'-C2'	-2.64	1.36	1.43
6	B	301	G4P	PC-O3'	2.54	1.67	1.59
6	B	301	G4P	O2'-C2'	-2.51	1.36	1.43
6	A	301	G4P	C5-C6	2.50	1.53	1.44
6	B	301	G4P	C6-N1	2.49	1.43	1.38
6	A	301	G4P	O6-C6	-2.47	1.18	1.23
6	B	301	G4P	C5-C6	2.45	1.53	1.44
6	B	301	G4P	O6-C6	-2.44	1.18	1.23
6	A	301	G4P	C6-N1	2.27	1.43	1.38
6	A	301	G4P	PC-O3'	2.26	1.66	1.59
6	B	301	G4P	O3'-C3'	2.10	1.51	1.44

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	301	G4P	C1'-N9-C4	-5.80	109.35	126.49
6	B	301	G4P	C1'-N9-C4	-5.56	110.06	126.49
6	B	301	G4P	C5-C4-N3	-5.38	119.83	128.39
6	A	301	G4P	C1'-N9-C8	5.20	141.49	126.73
6	A	301	G4P	C5-C4-N3	-5.14	120.21	128.39
6	B	301	G4P	C1'-N9-C8	5.02	140.98	126.73
6	B	301	G4P	C2-N3-C4	4.93	120.79	112.30
6	A	301	G4P	C2-N3-C4	4.42	119.92	112.30
6	B	301	G4P	N9-C4-N3	3.64	133.24	125.95
6	A	301	G4P	N9-C4-N3	3.41	132.78	125.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	301	G4P	N9-C8-N7	-3.25	107.37	113.40
6	B	301	G4P	N9-C8-N7	-3.16	107.53	113.40
6	B	301	G4P	C2-N1-C6	-3.03	119.62	125.11
6	A	301	G4P	C2-N1-C6	-3.00	119.67	125.11
6	B	301	G4P	O6-C6-C5	-3.00	118.61	126.53
6	B	301	G4P	C5-C6-N1	2.75	120.25	113.25
6	A	301	G4P	C5-C6-N1	2.66	120.02	113.25
6	A	301	G4P	O6-C6-C5	-2.61	119.63	126.53
6	B	301	G4P	N2-C2-N1	2.22	121.44	116.76
6	B	301	G4P	N1-C2-N3	-2.13	119.42	123.32
6	A	301	G4P	C8-N7-C5	2.13	108.06	104.26
6	B	301	G4P	C8-N7-C5	2.12	108.04	104.26
6	B	301	G4P	C3'-C2'-C1'	2.00	104.30	99.89

There are no chirality outliers.

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	301	GOL	O1-C1-C2-C3
3	C	301	GOL	C1-C2-C3-O3
3	C	302	GOL	C1-C2-C3-O3
3	C	303	GOL	O1-C1-C2-C3
3	C	303	GOL	C1-C2-C3-O3
3	A	302	GOL	O1-C1-C2-C3
3	A	302	GOL	C1-C2-C3-O3
6	A	301	G4P	C5'-O5'-PA-O1A
6	A	301	G4P	C5'-O5'-PA-O2A
6	A	301	G4P	O4'-C4'-C5'-O5'
6	B	301	G4P	PC-O3C-PD-O2D
6	B	301	G4P	PC-O3C-PD-O3D
4	C	305	1PE	OH4-C13-C23-OH3
4	B	302	1PE	OH5-C14-C24-OH4
4	C	305	1PE	OH6-C15-C25-OH5
4	B	302	1PE	OH2-C12-C22-OH3
6	A	301	G4P	C3'-C4'-C5'-O5'
3	C	301	GOL	O2-C2-C3-O3
3	C	302	GOL	O2-C2-C3-O3
3	C	303	GOL	O1-C1-C2-O2
3	C	303	GOL	O2-C2-C3-O3
3	A	302	GOL	O1-C1-C2-O2
3	A	302	GOL	O2-C2-C3-O3
6	B	301	G4P	O4'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
4	C	305	1PE	OH7-C16-C26-OH6
3	C	301	GOL	O1-C1-C2-O2
4	C	305	1PE	OH5-C14-C24-OH4
3	C	302	GOL	O1-C1-C2-O2
6	B	301	G4P	PB-O3A-PA-O5'
4	B	302	1PE	OH7-C16-C26-OH6
4	C	305	1PE	C14-C24-OH4-C13
4	B	302	1PE	OH6-C15-C25-OH5
6	A	301	G4P	C5'-O5'-PA-O3A
4	C	305	1PE	C13-C23-OH3-C22
3	C	304	GOL	O1-C1-C2-O2
6	A	301	G4P	PB-O3A-PA-O1A
6	B	301	G4P	PD-O3C-PC-O1C
4	B	302	1PE	OH4-C13-C23-OH3
4	C	305	1PE	C12-C22-OH3-C23
6	A	301	G4P	C3'-O3'-PC-O1C
6	A	301	G4P	C3'-O3'-PC-O2C
6	B	301	G4P	PB-O3A-PA-O2A
4	C	305	1PE	OH2-C12-C22-OH3
6	B	301	G4P	PD-O3C-PC-O2C

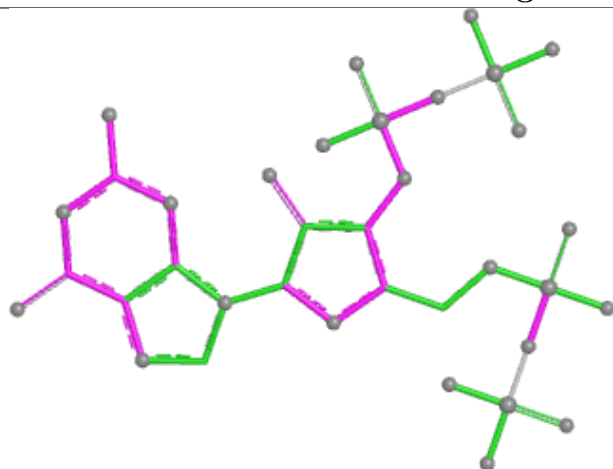
There are no ring outliers.

2 monomers are involved in 2 short contacts:

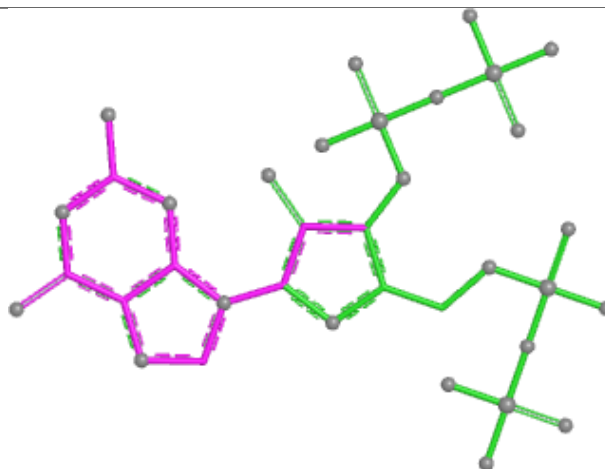
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	301	G4P	1	0
4	B	302	1PE	1	0

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.

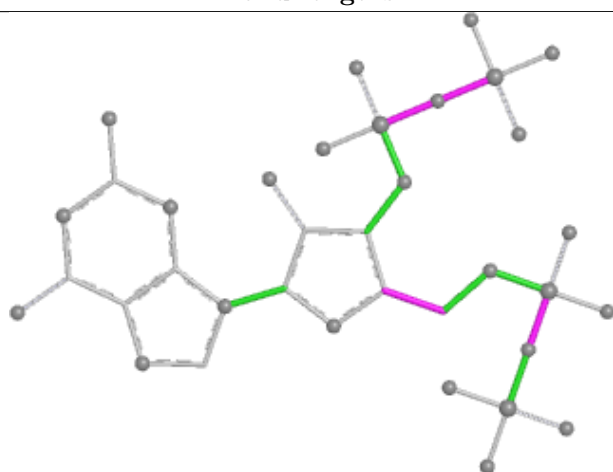
Ligand G4P B 301



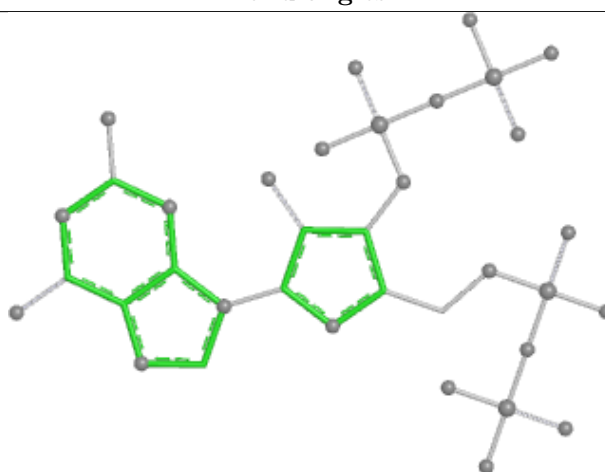
Bond lengths



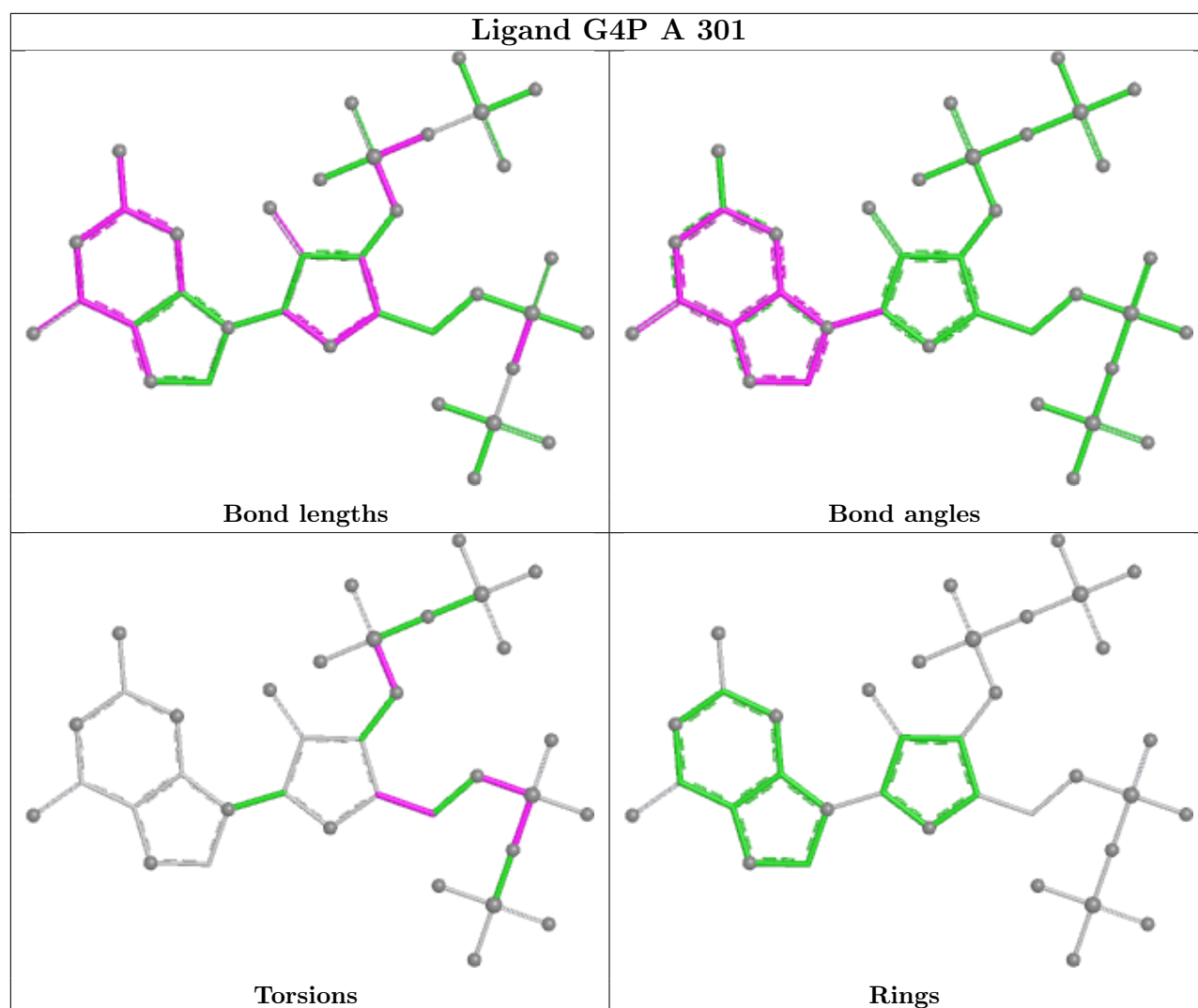
Bond angles



Torsions



Rings



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	C	195/211 (92%)	0.11	1 (0%) 87 82	37, 59, 88, 99	0
1	D	194/211 (91%)	0.22	1 (0%) 87 82	38, 65, 95, 103	0
2	A	203/204 (99%)	-0.13	2 (0%) 79 72	33, 53, 82, 99	0
2	B	201/204 (98%)	-0.15	3 (1%) 72 63	32, 50, 75, 94	0
All	All	793/830 (95%)	0.01	7 (0%) 81 74	32, 56, 88, 103	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	173	TYR	3.6
2	B	148	SER	2.7
1	D	130	LEU	2.5
2	A	0	ALA	2.3
2	B	0	ALA	2.1
2	A	41	GLU	2.1
2	B	151	ASP	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

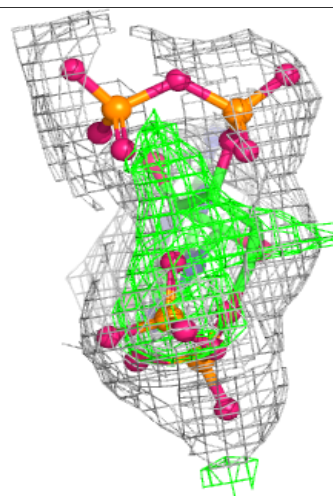
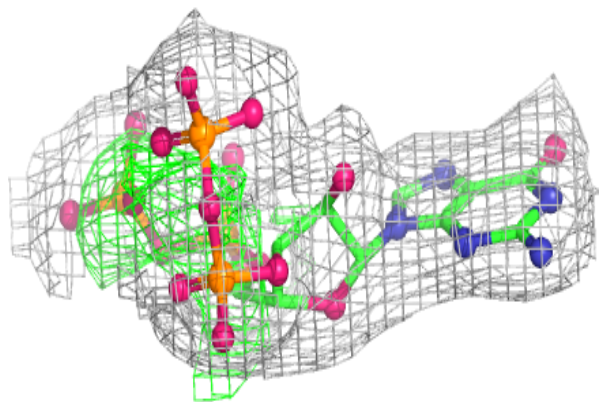
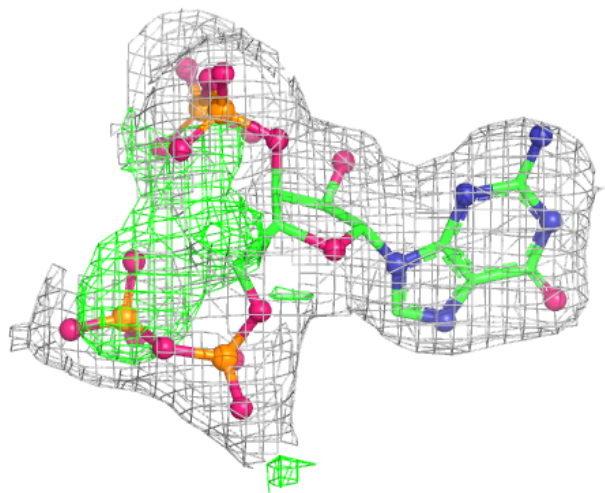
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	GOL	A	302	6/6	0.63	0.13	88,106,116,119	0
3	GOL	C	302	6/6	0.75	0.20	62,75,86,86	0
4	1PE	C	305	16/16	0.78	0.12	62,79,90,93	0
3	GOL	C	303	6/6	0.82	0.34	60,68,69,74	0
3	GOL	C	304	6/6	0.85	0.13	69,83,96,96	0
5	MG	B	303	1/1	0.85	0.22	60,60,60,60	0
3	GOL	C	301	6/6	0.87	0.16	60,72,88,88	0
4	1PE	B	302	16/16	0.88	0.10	55,70,86,91	0
6	G4P	A	301	36/36	0.90	0.10	32,44,65,163	0
6	G4P	B	301	36/36	0.93	0.08	32,47,58,78	0
5	MG	C	306	1/1	0.96	0.14	50,50,50,50	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.

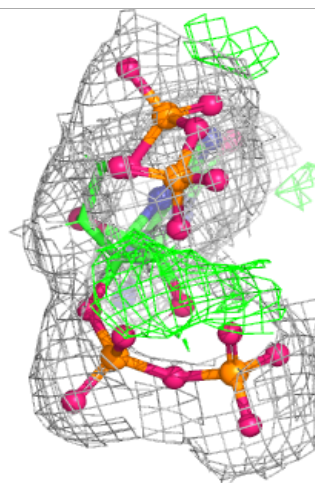
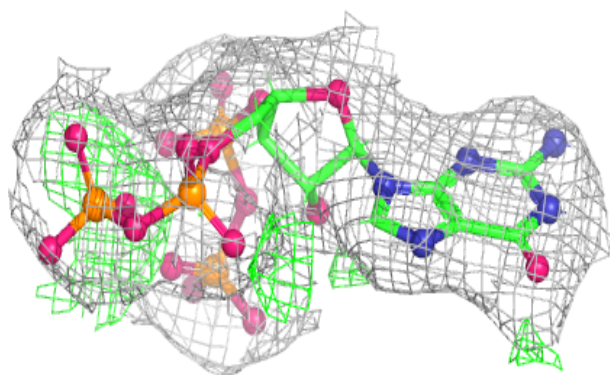
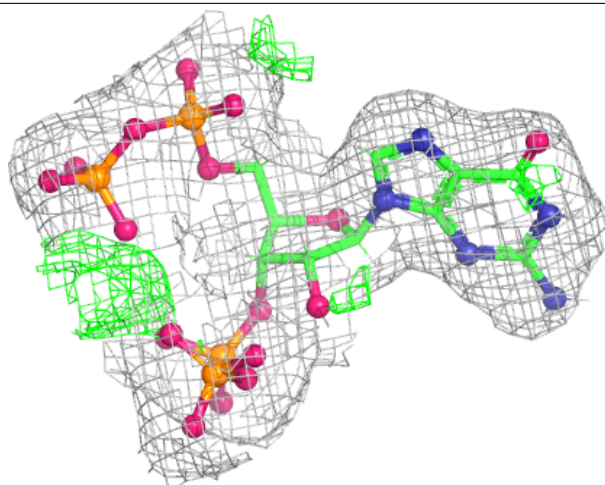
Electron density around G4P A 301:

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



Electron density around G4P B 301:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.