



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:52 AM UTC

PDB ID : 5D7J / pdb_00005d7j
Title : Structure of human MR1-5-OP-RU in complex with human MAIT
M33.64(Y95alphaF) TCR
Authors : Keller, A.N.; Woolley, R.E.; Rossjohn, J.
Deposited on : 2015-08-14
Resolution : 1.97 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:

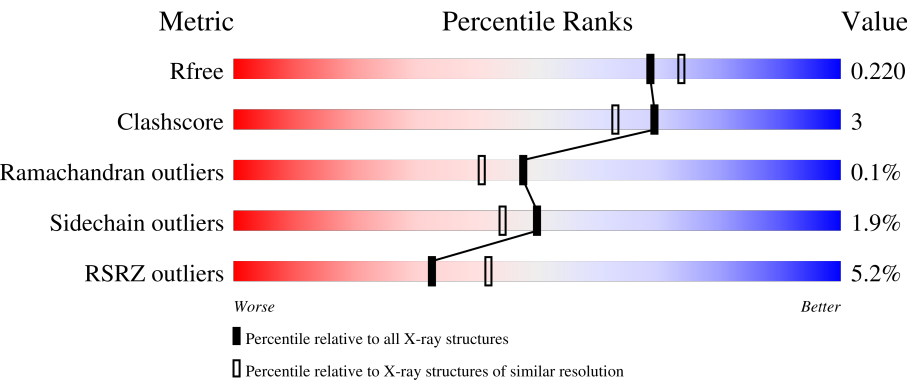
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.97 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.



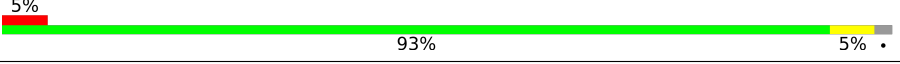
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	1506 (1.98-1.98)
Clashscore	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)
RSRZ outliers	180081	1506 (1.98-1.98)

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	204	4% 91% 7%
1	G	204	5% 85% 11%
2	B	245	3% 90% 7%
2	H	245	3% 87% 11%
3	C	271	4% 90% 6%

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Mol	Chain	Length	Quality of chain
3	E	271	 5% 93% 5% •
4	D	100	 4% 85% 12% ••
4	F	100	 20% 87% 9% ••

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 14052 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 M33.64 TCR Alpha Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	200	Total	C	N	O	S	0	1	0
			1542	976	244	312	10			
1	G	196	Total	C	N	O	S	0	2	0
			1521	961	244	306	10			

- Molecule 2 is a protein called M33.64 TCR Beta Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	240	Total	C	N	O	S	0	1	0
			1852	1162	320	362	8			
2	H	242	Total	C	N	O	S	0	1	0
			1873	1176	332	357	8			

- Molecule 3 is a protein called Major histocompatibility complex class I-related gene protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	263	Total	C	N	O	S	0	2	0
			2162	1386	376	390	10			
3	E	265	Total	C	N	O	S	0	5	0
			2175	1390	379	395	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	MET	-	initiating methionine	UNP Q95460
C	261	SER	CYS	conflict	UNP Q95460
E	0	MET	-	initiating methionine	UNP Q95460
E	261	SER	CYS	conflict	UNP Q95460

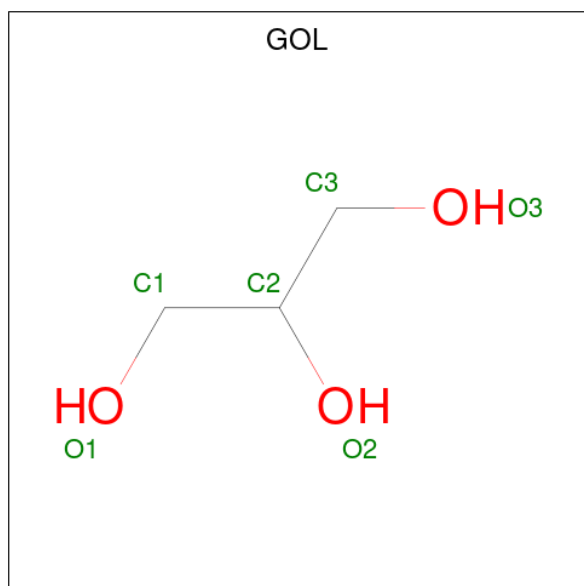
- Molecule 4 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	98	Total	C	N	O	S	0	0	0
			816	520	138	156	2			
4	F	98	Total	C	N	O	S	0	0	0
			783	504	134	142	3			

There are 2 discrepancies between the modelled and reference sequences:

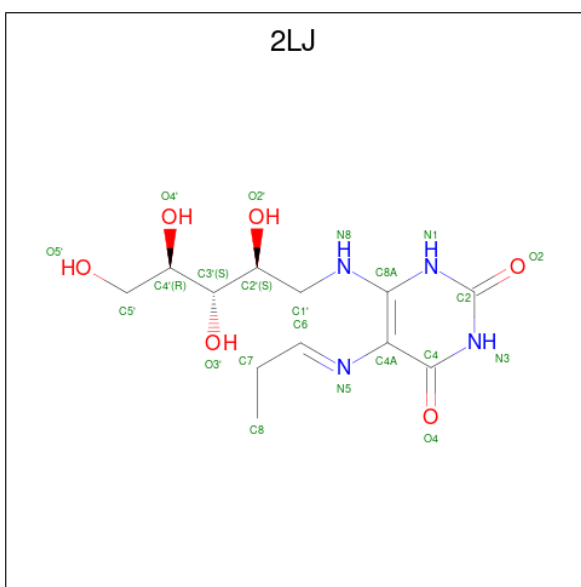
Chain	Residue	Modelled	Actual	Comment	Reference
D	0	MET	-	initiating methionine	UNP P61769
F	0	MET	-	initiating methionine	UNP P61769

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			6	3	3		
5	D	1	Total	C	O	0	0
			6	3	3		
5	H	1	Total	C	O	0	0
			6	3	3		
5	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is 1-deoxy-1-({2,6-dioxo-5-[(E)-propylideneamino]-1,2,3,6-tetrahydropyrimidin-4-yl}amino)-D-ribitol (CCD ID: 2LJ) (formula: $C_{12}H_{20}N_4O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	C	1	Total	C	N	O	0	0
			22	12	4	6		
6	E	1	Total	C	N	O	0	0
			22	12	4	6		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	145	Total	O	0	0
			145	145		
7	B	135	Total	O	0	0
			135	135		
7	C	258	Total	O	0	0
			258	258		
7	D	91	Total	O	0	0
			91	91		
7	E	185	Total	O	0	0
			185	185		
7	F	43	Total	O	0	0
			43	43		
7	G	156	Total	O	0	0
			156	156		
7	H	247	Total	O	0	0
			247	247		

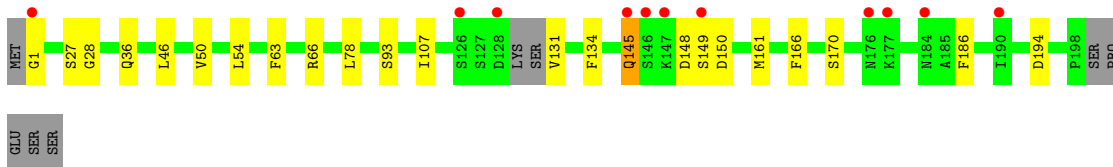
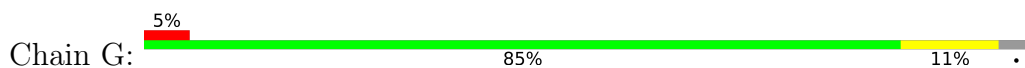
3 Residue-property plots [i](#)

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.

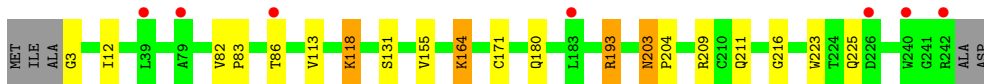
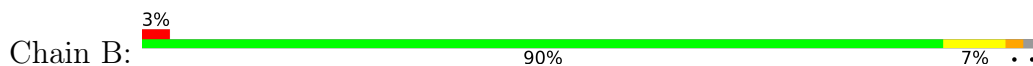
- Molecule 1: M33.64 TCR Alpha Chain



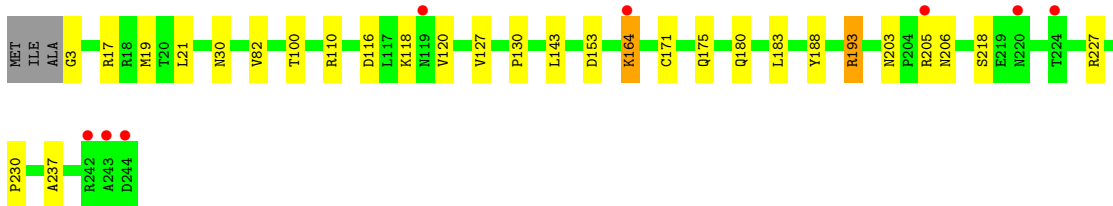
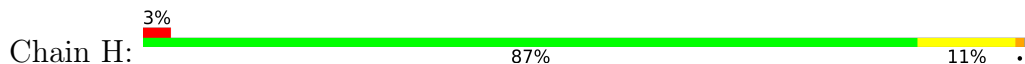
- Molecule 1: M33.64 TCR Alpha Chain



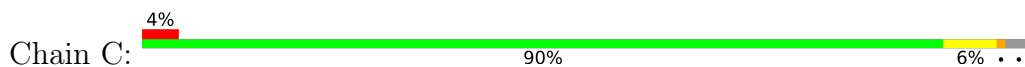
- Molecule 2: M33.64 TCR Beta Chain



- Molecule 2: M33.64 TCR Beta Chain



- Molecule 3: Major histocompatibility complex class I-related gene protein

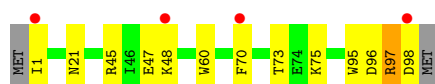
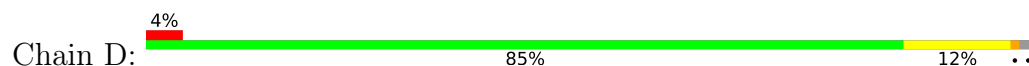




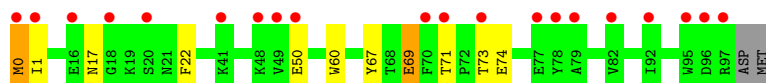
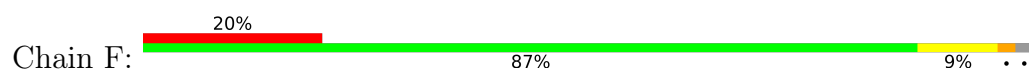
- Molecule 3: Major histocompatibility complex class I-related gene protein



- Molecule 4: Beta-2-microglobulin



- Molecule 4: Beta-2-microglobulin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	214.60Å 69.59Å 141.47Å 90.00° 103.81° 90.00°	Depositor
Resolution (Å)	35.76 – 1.97 35.76 – 1.97	Depositor EDS
% Data completeness (in resolution range)	93.7 (35.76-1.97) 93.9 (35.76-1.97)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 1.97Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.179 , 0.222 0.179 , 0.220	Depositor DCC
R_{free} test set	6828 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	22.2	Xtriage
Anisotropy	0.580	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.5	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	0.95	EDS
Total number of atoms	14052	wwPDB-VP
Average B, all atoms (Å ²)	32.0	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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 2LJ, 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	A	0.49	0/1580	0.79	1/2146 (0.0%)
1	G	0.49	0/1557	0.77	0/2113
2	B	0.44	0/1900	0.77	4/2594 (0.2%)
2	H	0.47	0/1924	0.77	0/2622
3	C	0.54	1/2235 (0.0%)	0.76	0/3039
3	E	0.44	0/2252	0.73	0/3063
4	D	0.47	0/839	0.75	0/1138
4	F	0.43	0/806	0.74	0/1097
All	All	0.48	1/13093 (0.0%)	0.76	5/17812 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	152	TYR	C-O	-6.85	1.16	1.24

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	25	GLN	N-CA-C	5.55	118.07	107.75
2	B	203	ASN	CA-C-N	5.24	125.04	119.28
2	B	203	ASN	C-N-CA	5.24	125.04	119.28
2	B	180	GLN	CA-C-N	5.12	125.40	119.47
2	B	180	GLN	C-N-CA	5.12	125.40	119.47

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	1542	0	1437	11	0
1	G	1521	0	1414	13	0
2	B	1852	0	1732	14	0
2	H	1873	0	1773	18	0
3	C	2162	0	2036	15	0
3	E	2175	0	2036	8	0
4	D	816	0	774	7	0
4	F	783	0	735	6	0
5	B	6	0	8	1	0
5	D	6	0	8	0	0
5	H	12	0	16	0	0
6	C	22	0	18	1	0
6	E	22	0	18	1	0
7	A	145	0	0	2	0
7	B	135	0	0	2	0
7	C	258	0	0	7	0
7	D	91	0	0	1	0
7	E	185	0	0	0	0
7	F	43	0	0	0	0
7	G	156	0	0	1	0
7	H	247	0	0	6	0
All	All	14052	0	12005	86	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 3.

All (86) 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:B:86:THR:HG22	2:B:113:VAL:H	1.26	0.97
2:B:118:LYS:NZ	2:B:225:GLN:OE1	2.19	0.75
1:A:1:GLY:HA2	1:A:27:SER:H	1.54	0.73
1:A:170:SER:OG	2:B:193:ARG:NE	2.22	0.72
4:D:45:ARG:NH1	4:D:47:GLU:OE2	2.25	0.70
2:B:83:PRO:O	2:B:86:THR:HG23	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:1:GLY:HA2	1:G:27:SER:H	1.56	0.69
1:A:170:SER:O	2:B:193:ARG:NH1	2.27	0.66
1:A:19:GLN:NE2	7:A:301:HOH:O	2.26	0.65
4:F:0:MET:HG2	4:F:1:ILE:N	2.14	0.62
2:H:110:ARG:NH2	2:H:153:ASP:OD1	2.27	0.62
3:C:46:ARG:NH2	7:C:705:HOH:O	2.33	0.61
4:F:22:PHE:CZ	4:F:69:GLU:HG3	2.36	0.61
1:G:54:LEU:HD23	1:G:63:PHE:HB2	1.83	0.60
2:H:17[A]:ARG:NH2	7:H:407:HOH:O	2.35	0.60
3:C:1:ARG:N	7:C:706:HOH:O	2.34	0.59
2:B:86:THR:HG22	2:B:113:VAL:N	2.09	0.59
3:C:152:TYR:OH	7:C:701:HOH:O	2.10	0.59
4:D:73:THR:O	4:D:97:ARG:NH2	2.36	0.59
3:E:6[A]:ARG:NH2	3:E:29:ASP:O	2.30	0.59
3:E:113:ALA:HB2	4:F:60:TRP:CE2	2.38	0.59
1:G:145:GLN:NE2	7:G:304:HOH:O	2.37	0.57
1:A:125:LYS:O	3:C:130:LEU:HD23	2.05	0.57
1:A:157:CYS:HB3	2:B:193:ARG:NH1	2.21	0.56
2:H:3:GLY:N	7:H:409:HOH:O	2.40	0.55
1:A:159:LEU:HB3	2:B:171:CYS:HB2	1.89	0.55
4:F:17:ASN:HD21	4:F:74:GLU:N	2.05	0.55
2:H:171:CYS:O	2:H:193:ARG:NH2	2.39	0.55
2:H:180:GLN:HB3	2:H:183:LEU:HD13	1.89	0.54
2:H:116:ASP:O	7:H:402:HOH:O	2.18	0.54
3:C:113:ALA:HB2	4:D:60:TRP:CE2	2.44	0.53
1:G:28:GLY:HA3	1:G:93[B]:SER:OG	2.08	0.53
1:G:161:MET:HE2	1:G:166:PHE:CD1	2.44	0.52
1:G:194:ASP:OD1	1:G:194:ASP:N	2.33	0.52
3:C:160:GLU:OE2	7:C:702:HOH:O	2.19	0.52
4:D:48:LYS:HE3	7:D:247:HOH:O	2.08	0.52
3:C:255:SER:HB3	3:C:268:GLN:OE1	2.10	0.51
3:E:215:MET:HG3	3:E:257:HIS:CD2	2.46	0.51
2:B:209:ARG:NH1	2:B:211:GLN:HB2	2.25	0.51
1:G:170:SER:OG	2:H:193:ARG:HD3	2.10	0.51
2:B:223:TRP:CZ2	2:B:225:GLN:HB2	2.46	0.51
3:C:4:SER:HB3	3:C:99:GLU:HG2	1.93	0.50
2:H:130:PRO:HD3	2:H:143:LEU:HG	1.92	0.50
4:D:21:ASN:HB3	4:D:70:PHE:CE2	2.47	0.49
1:G:148:ASP:OD1	1:G:150:ASP:N	2.33	0.49
1:G:36:GLN:HB2	1:G:46:LEU:HD11	1.94	0.49
6:C:600:2LJ:O4	6:C:600:2LJ:H1	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:205:ARG:NH2	7:H:401:HOH:O	2.05	0.48
3:C:1:ARG:N	3:C:102:GLU:HG3	2.28	0.48
7:C:726:HOH:O	2:H:100:THR:HG23	2.13	0.48
2:H:164:LYS:HG3	7:H:450:HOH:O	2.13	0.48
1:G:134:PHE:HB2	1:G:186:PHE:CE2	2.50	0.47
3:E:215:MET:HE2	3:E:257:HIS:CE1	2.49	0.47
1:A:51:LEU:HD12	3:E:151:LEU:HD22	1.97	0.47
1:A:79:GLN:NE2	7:A:302:HOH:O	2.28	0.47
1:G:148:ASP:OD1	1:G:149:SER:N	2.47	0.46
3:E:152:TYR:CD2	3:E:152:TYR:C	2.91	0.46
6:E:600:2LJ:O4	6:E:600:2LJ:H1	2.16	0.46
2:B:3:GLY:N	7:B:409:HOH:O	2.47	0.46
4:D:96:ASP:O	4:D:98:ASP:N	2.49	0.46
3:C:185:ARG:NH1	7:C:704:HOH:O	2.32	0.45
3:C:77:LEU:HD13	3:C:92:TYR:HB2	1.98	0.45
2:H:118:LYS:O	2:H:227:ARG:NH2	2.49	0.45
5:B:301:GOL:O2	7:B:401:HOH:O	2.21	0.45
2:H:203:ASN:HB3	2:H:206:ASN:ND2	2.33	0.44
2:B:12:ILE:HD12	2:B:216:GLY:HA2	2.00	0.44
3:E:77:LEU:HD13	3:E:92:TYR:HB2	2.01	0.43
2:H:19:MET:HE3	2:H:21:LEU:HD23	2.01	0.43
4:D:95:TRP:CZ2	4:D:97:ARG:HA	2.54	0.43
2:H:120:VAL:HG12	2:H:230:PRO:HB2	2.00	0.43
3:C:185:ARG:NH2	3:C:201:LYS:HD3	2.34	0.43
1:G:78:LEU:HD22	1:G:107:ILE:HD11	2.00	0.43
3:C:156:TRP:CZ3	3:C:161:CYS:HB2	2.54	0.42
1:G:50:VAL:O	1:G:66:ARG:HD3	2.20	0.42
3:C:1:ARG:HD3	7:C:916:HOH:O	2.20	0.42
2:H:30:ASN:ND2	7:H:412:HOH:O	2.41	0.42
2:H:127:VAL:HG23	2:H:237:ALA:HB3	2.00	0.42
2:H:153:ASP:HB3	2:H:188:TYR:CE2	2.55	0.41
4:F:50:GLU:HB3	4:F:67:TYR:CZ	2.56	0.41
2:B:164:LYS:HD3	2:B:164:LYS:HA	1.74	0.41
1:A:61[B]:SER:OG	1:A:74:LEU:HB3	2.21	0.41
1:A:177:LYS:HA	1:A:177:LYS:HD3	1.83	0.41
2:B:203:ASN:HA	2:B:204:PRO:HD3	1.88	0.41
3:E:156:TRP:CZ3	3:E:161:CYS:HB2	2.56	0.41
3:C:16:ILE:H	3:C:16:ILE:HG13	1.59	0.41
4:F:17:ASN:OD1	4:F:73:THR:HA	2.20	0.41

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	199/204 (98%)	193 (97%)	6 (3%)	0	100	100
1	G	194/204 (95%)	189 (97%)	5 (3%)	0	100	100
2	B	239/245 (98%)	233 (98%)	6 (2%)	0	100	100
2	H	241/245 (98%)	235 (98%)	6 (2%)	0	100	100
3	C	261/271 (96%)	254 (97%)	7 (3%)	0	100	100
3	E	265/271 (98%)	259 (98%)	6 (2%)	0	100	100
4	D	96/100 (96%)	95 (99%)	0	1 (1%)	12	5
4	F	96/100 (96%)	95 (99%)	1 (1%)	0	100	100
All	All	1591/1640 (97%)	1553 (98%)	37 (2%)	1 (0%)	48	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	97	ARG

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	170/181 (94%)	169 (99%)	1 (1%)	78	76
1	G	167/181 (92%)	165 (99%)	2 (1%)	63	58
2	B	197/205 (96%)	191 (97%)	6 (3%)	36	27
2	H	197/205 (96%)	192 (98%)	5 (2%)	42	33

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	226/241 (94%)	223 (99%)	3 (1%)	61	57
3	E	226/241 (94%)	222 (98%)	4 (2%)	51	46
4	D	92/95 (97%)	90 (98%)	2 (2%)	45	38
4	F	84/95 (88%)	81 (96%)	3 (4%)	31	20
All	All	1359/1444 (94%)	1333 (98%)	26 (2%)	50	44

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	154	THR
2	B	82	VAL
2	B	118	LYS
2	B	131	SER
2	B	155	VAL
2	B	164	LYS
2	B	193	ARG
3	C	46	ARG
3	C	221	ILE
3	C	268	GLN
4	D	1	ILE
4	D	75	LYS
3	E	1[A]	ARG
3	E	1[B]	ARG
3	E	25	VAL
3	E	102	GLU
4	F	0	MET
4	F	69	GLU
4	F	71	THR
1	G	131	VAL
1	G	145	GLN
2	H	82	VAL
2	H	164	LYS
2	H	175	GLN
2	H	193	ARG
2	H	218	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	ASN

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Mol	Chain	Res	Type
1	A	112	GLN
1	A	113	ASN
2	B	11	GLN
3	C	42	GLN
3	C	134	ASN
3	C	177	GLN
3	C	187	ASN
3	E	64	GLN
3	E	177	GLN
1	G	37	GLN
1	G	96	GLN
1	G	140	GLN
1	G	145	GLN
2	H	30	ASN
2	H	37	GLN

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
5	GOL	H	301	-	5,5,5	0.48	0	5,5,5	0.63	0
5	GOL	H	302	-	5,5,5	0.26	0	5,5,5	1.34	1 (20%)
5	GOL	B	301	-	5,5,5	0.44	0	5,5,5	0.54	0
6	2LJ	E	600	3	22,22,22	0.66	1 (4%)	24,29,29	1.22	4 (16%)
6	2LJ	C	600	3	22,22,22	0.67	1 (4%)	24,29,29	1.13	3 (12%)
5	GOL	D	101	-	5,5,5	0.45	0	5,5,5	0.23	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
5	GOL	H	301	-	-	0/4/4/4	-
5	GOL	H	302	-	-	2/4/4/4	-
5	GOL	B	301	-	-	0/4/4/4	-
6	2LJ	E	600	3	-	1/18/19/19	0/1/1/1
6	2LJ	C	600	3	-	1/18/19/19	0/1/1/1
5	GOL	D	101	-	-	0/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	E	600	2LJ	C7-C6	-2.50	1.46	1.49
6	C	600	2LJ	C7-C6	-2.27	1.47	1.49

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	E	600	2LJ	C4'-C3'-C2'	3.03	118.61	113.57
5	H	302	GOL	C3-C2-C1	-2.86	101.31	111.80
6	C	600	2LJ	C4A-C8A-N1	2.64	121.31	115.89
6	E	600	2LJ	C2'-C1'-N8	2.44	118.29	111.53
6	E	600	2LJ	C8-C7-C6	2.31	117.84	113.75
6	C	600	2LJ	C4'-C3'-C2'	2.13	117.12	113.57
6	E	600	2LJ	C4A-C8A-N1	2.09	120.19	115.89
6	C	600	2LJ	C2'-C1'-N8	2.06	117.23	111.53

There are no chirality outliers.

All (4) torsion outliers are listed below:

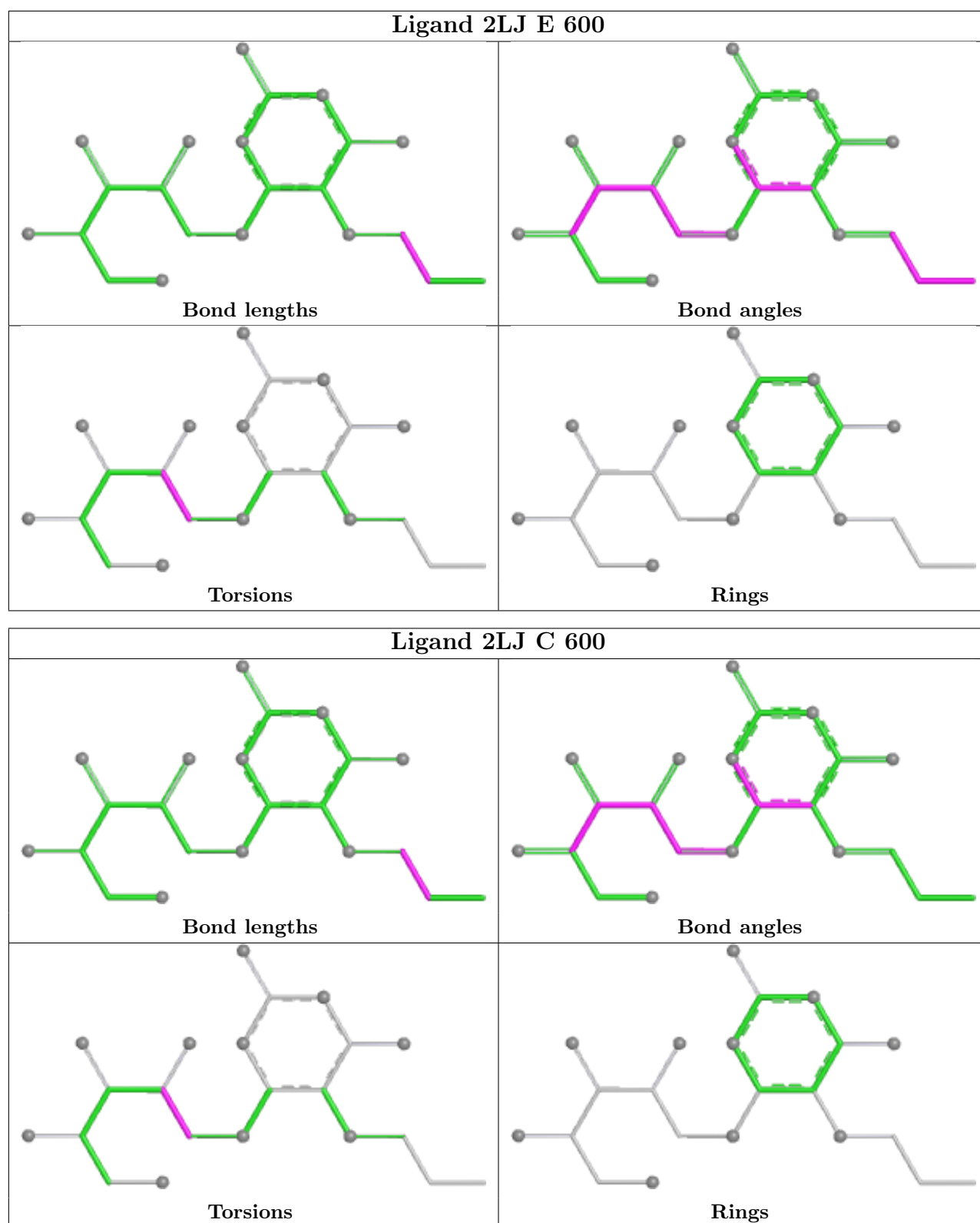
Mol	Chain	Res	Type	Atoms
5	H	302	GOL	C1-C2-C3-O3
5	H	302	GOL	O2-C2-C3-O3
6	E	600	2LJ	N8-C1'-C2'-C3'
6	C	600	2LJ	N8-C1'-C2'-C3'

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	301	GOL	1	0
6	E	600	2LJ	1	0
6	C	600	2LJ	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	200/204 (98%)	0.11	8 (4%)	42	52	14, 29, 61, 72	1 (0%)
1	G	196/204 (96%)	0.11	11 (5%)	30	38	7, 28, 73, 85	2 (1%)
2	B	240/245 (97%)	0.33	7 (2%)	53	63	15, 36, 57, 82	1 (0%)
2	H	242/245 (98%)	-0.13	8 (3%)	49	59	12, 25, 51, 77	1 (0%)
3	C	263/271 (97%)	-0.18	11 (4%)	40	50	7, 21, 56, 86	2 (0%)
3	E	265/271 (97%)	0.12	14 (5%)	32	41	12, 30, 59, 81	5 (1%)
4	D	98/100 (98%)	-0.03	4 (4%)	41	51	14, 29, 58, 70	0
4	F	98/100 (98%)	0.95	20 (20%)	3	3	24, 54, 81, 92	0
All	All	1602/1640 (97%)	0.10	83 (5%)	33	42	7, 29, 65, 92	12 (0%)

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	224	THR	4.7
3	C	221	ILE	4.6
1	G	1	GLY	4.5
3	C	246	LEU	4.4
1	G	128	ASP	4.0
1	A	1	GLY	3.9
1	G	176	ASN	3.5
2	H	243	ALA	3.5
3	C	195	VAL	3.5
3	C	17	HIS	3.5
2	B	242	ARG	3.5
4	F	18	GLY	3.5
1	G	126	SER	3.4
3	C	222	VAL	3.3
3	C	211	TYR	3.3
2	B	183	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
4	F	41	LYS	3.2
4	D	98	ASP	3.2
2	B	226	ASP	3.1
1	G	147	LYS	3.1
2	H	220	ASN	3.1
4	D	48	LYS	3.1
4	F	73	THR	3.1
2	H	205	ARG	3.0
1	G	184	ASN	3.0
3	E	211	TYR	3.0
4	F	82	VAL	3.0
3	E	188	ARG	3.0
3	E	17	HIS	2.9
4	F	50	GLU	2.9
1	A	163	SER	2.9
2	H	244	ASP	2.9
3	E	195	VAL	2.8
1	G	145	GLN	2.8
3	C	218	GLY	2.8
1	G	177	LYS	2.8
4	F	71	THR	2.8
1	A	200	PRO	2.7
4	F	49	VAL	2.7
3	C	18	GLY	2.6
4	F	97	ARG	2.6
4	F	48	LYS	2.6
1	A	145	GLN	2.6
4	F	95	TRP	2.6
4	F	0	MET	2.6
1	A	2	GLN	2.5
3	E	249	GLN	2.5
3	E	194	GLY	2.5
2	H	242	ARG	2.5
4	F	79	ALA	2.5
3	E	16	ILE	2.4
1	G	146	SER	2.4
4	D	70	PHE	2.4
4	D	1	ILE	2.4
2	B	79	ALA	2.4
1	G	149	SER	2.4
2	H	164	LYS	2.4
3	E	196	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	154	THR	2.3
4	F	77	GLU	2.3
1	A	70	TYR	2.3
3	C	16	ILE	2.3
3	E	218	GLY	2.3
3	E	52	GLU	2.3
3	E	223	GLN	2.3
2	B	86	THR	2.3
3	C	245	GLU	2.2
4	F	70	PHE	2.2
4	F	1	ILE	2.2
2	B	240	TRP	2.2
4	F	96	ASP	2.2
3	C	220	GLU	2.1
3	E	19	VAL	2.1
4	F	78	TYR	2.1
3	E	152	TYR	2.1
4	F	20	SER	2.1
2	H	119	ASN	2.1
3	E	221	ILE	2.0
4	F	16	GLU	2.0
1	A	184	ASN	2.0
2	B	39	LEU	2.0
1	G	190	ILE	2.0
4	F	92	ILE	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 ⓘ

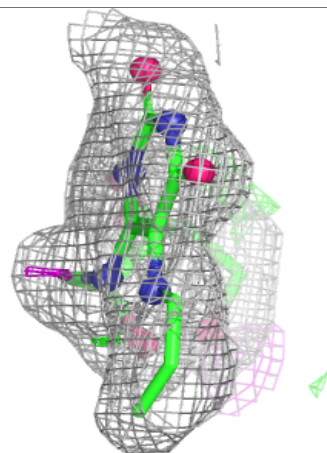
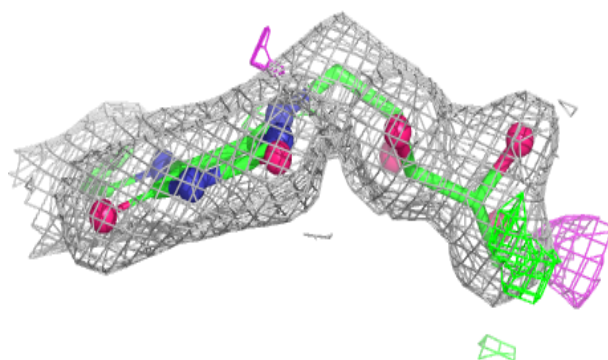
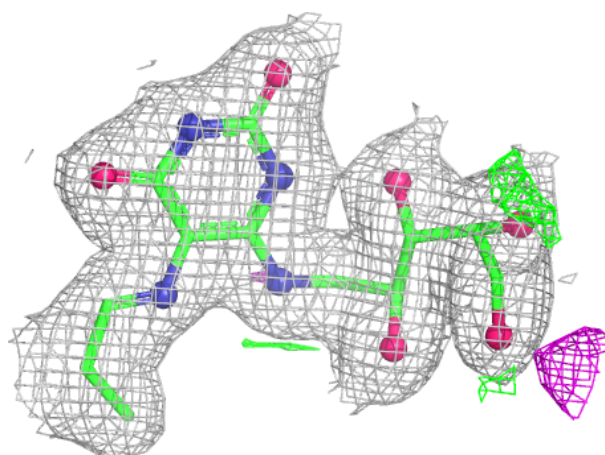
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	GOL	H	302	6/6	0.88	0.14	18,27,29,32	0
5	GOL	D	101	6/6	0.96	0.07	24,25,29,29	0
6	2LJ	E	600	22/22	0.96	0.06	12,15,24,27	0
6	2LJ	C	600	22/22	0.97	0.05	8,13,24,30	0
5	GOL	H	301	6/6	0.98	0.04	12,13,16,18	0
5	GOL	B	301	6/6	0.98	0.04	17,22,23,28	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.

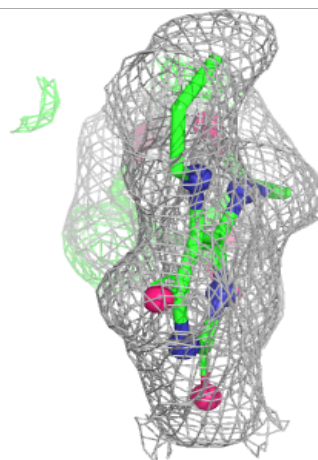
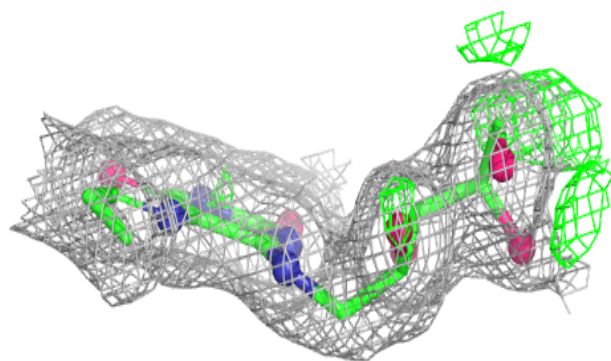
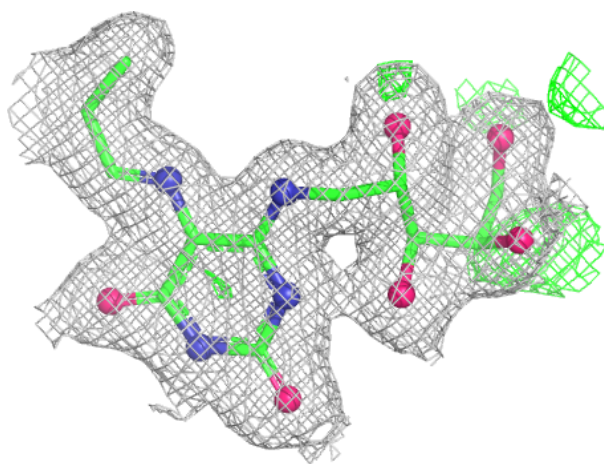
Electron density around 2LJ E 600:

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 2LJ C 600:

$2mF_o-DF_c$ (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.