



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 11:57 AM UTC

PDB ID : 6PVC / pdb_00006pvc
Title : Structure of human MAIT A-F7 TCR in complex with human MR1-DB28
Authors : Awad, W.; Rossjohn, J.
Deposited on : 2019-07-20
Resolution : 1.96 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

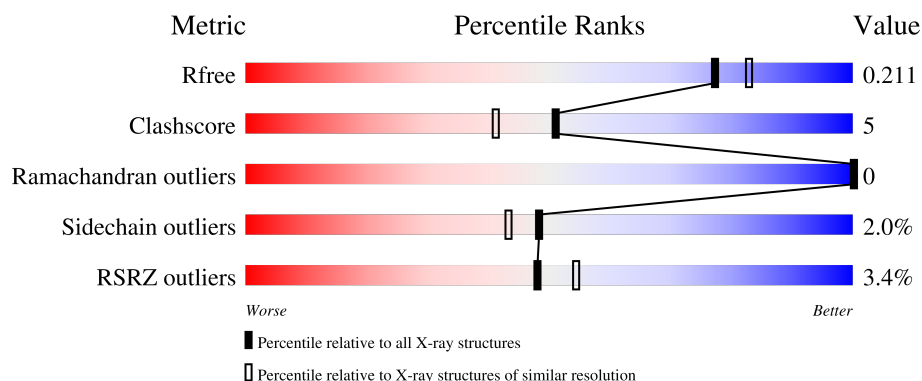
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.96 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	271	2% 91% 7% .
1	C	271	3% 84% 14% .
2	B	100	4% 94% ...
2	D	100	% 95% 5%
3	E	204	83% 10% 6%

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Mol	Chain	Length	Quality of chain
3	G	204	% 91% 7%
4	F	246	11% 78% 20%
4	H	246	2% 88% 11%

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 14903 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 Major histocompatibility complex class I-related gene protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	10	0
			2216	1419	383	402	12			
1	C	267	Total	C	N	O	S	0	17	0
			2288	1475	389	412	12			

There are 4 discrepancies between the modelled and reference sequences:

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	98	Total	C	N	O	S	0	1	0
			782	502	134	143	3			
2	D	100	Total	C	N	O	S	0	3	0
			831	533	138	156	4			

There are 2 discrepancies between the modelled and reference sequences:

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

- Molecule 3 is a protein called Human TCR alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	192	Total	C	N	O	S	0	10	0
			1540	979	244	306	11			

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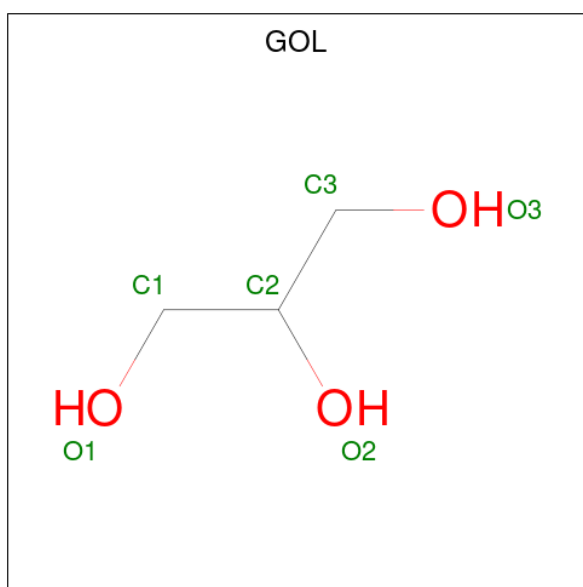
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	201	Total	C	N	O	S	0	14	0
			1633	1036	259	328	10			

- Molecule 4 is a protein called Human TCR beta chain.

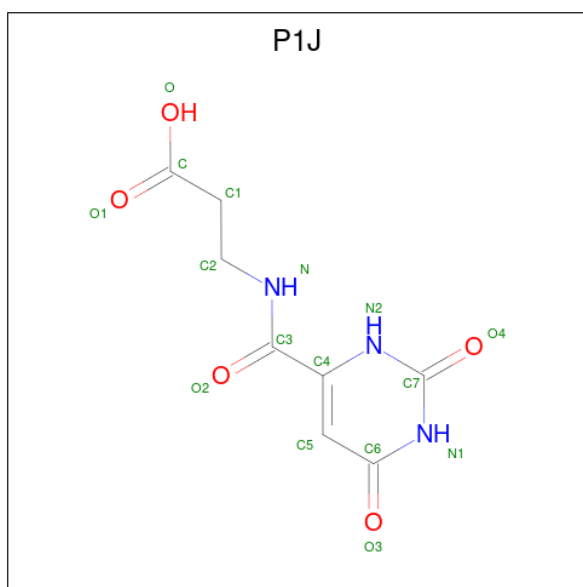
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	242	Total	C	N	O	S	0	21	0
			1989	1255	340	380	14			
4	H	242	Total	C	N	O	S	0	13	0
			1955	1239	334	368	14			

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



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

- Molecule 6 is N-(2,6-dioxo-1,2,3,6-tetrahydropyrimidine-4-carbonyl)-beta-alanine (CCD ID: P1J) (formula: $C_8H_9N_3O_5$) (labeled as "Ligand of Interest" by depositor).

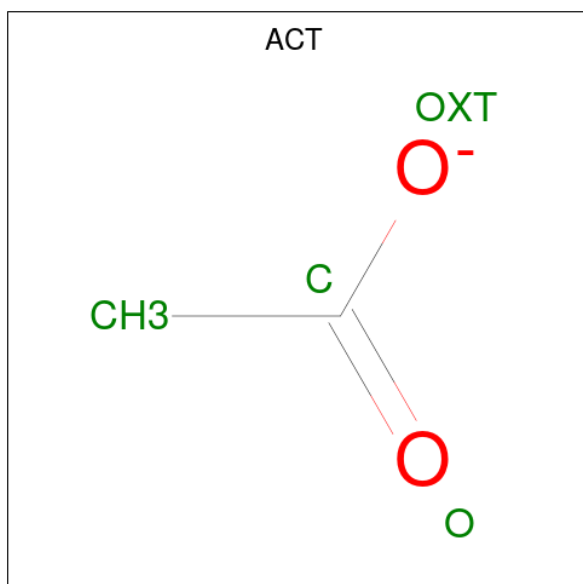


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			16	8	3	5		
6	C	1	Total	C	N	O	0	0
			16	8	3	5		

- Molecule 7 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	D	1	Total	Na	0	0
			1	1		

- Molecule 8 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	H	1	Total	C	O	0	0
			4	2	2		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	237	Total	O	0	0
			237	237		
9	B	73	Total	O	0	0
			73	73		
9	C	334	Total	O	0	0
			334	334		
9	D	117	Total	O	0	0
			117	117		
9	E	168	Total	O	0	0
			168	168		
9	F	207	Total	O	0	0
			207	207		
9	G	243	Total	O	0	0
			243	243		
9	H	241	Total	O	0	0
			241	241		

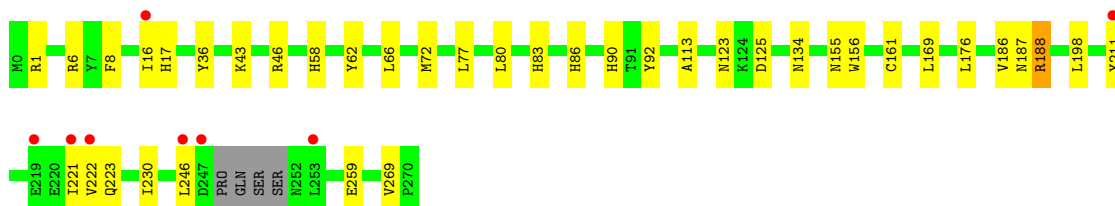
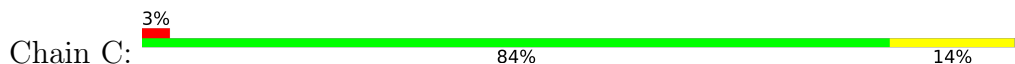
3 Residue-property plots [i](#)

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: Major histocompatibility complex class I-related gene protein



- Molecule 1: Major histocompatibility complex class I-related gene protein



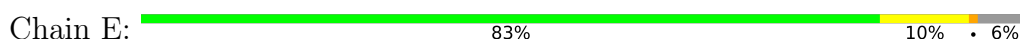
- Molecule 2: Beta-2-microglobulin

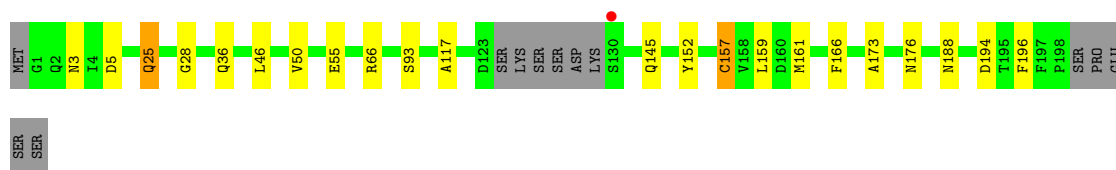


- Molecule 2: Beta-2-microglobulin



- Molecule 3: Human TCR alpha chain

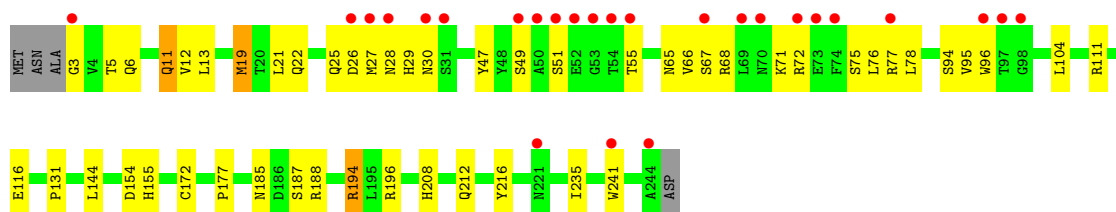
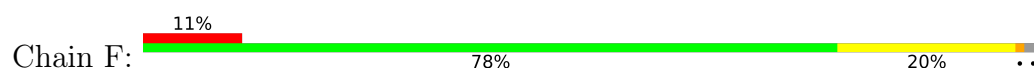




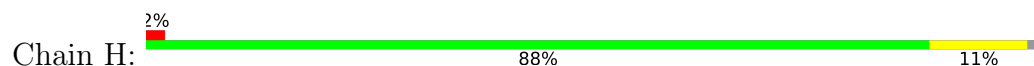
• Molecule 3: Human TCR alpha chain



• Molecule 4: Human TCR beta chain



• Molecule 4: Human TCR beta chain



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	214.29Å 69.83Å 141.62Å 90.00° 103.32° 90.00°	Depositor
Resolution (Å)	45.42 – 1.96 45.42 – 1.96	Depositor EDS
% Data completeness (in resolution range)	99.8 (45.42-1.96) 99.9 (45.42-1.96)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 1.97Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.169 , 0.212 0.170 , 0.211	Depositor DCC
R_{free} test set	7448 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	35.5	Xtriage
Anisotropy	0.190	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 59.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	14903	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.50% 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: NA, ACT, GOL, P1J

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	0.40	0/2303	0.55	0/3131
1	C	0.49	0/2405	0.62	0/3269
2	B	0.29	0/805	0.49	0/1099
2	D	0.43	0/863	0.56	0/1172
3	E	0.38	0/1601	0.56	0/2173
3	G	0.45	0/1708	0.63	0/2316
4	F	0.36	0/2093	0.56	0/2847
4	H	0.40	0/2052	0.59	0/2790
All	All	0.41	0/13830	0.58	0/18797

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	A	2216	0	2092	17	0
1	C	2288	0	2207	32	0
2	B	782	0	717	3	0
2	D	831	0	785	2	0
3	E	1540	0	1449	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	1633	0	1568	11	0
4	F	1989	0	1894	42	0
4	H	1955	0	1863	17	0
5	A	6	0	6	1	0
5	D	6	0	8	1	0
6	A	16	0	0	0	0
6	C	16	0	0	0	0
7	D	1	0	0	0	0
8	H	4	0	3	0	0
9	A	237	0	0	3	0
9	B	73	0	0	1	0
9	C	334	0	0	14	1
9	D	117	0	0	1	0
9	E	168	0	0	4	0
9	F	207	0	0	13	1
9	G	243	0	0	2	0
9	H	241	0	0	3	0
All	All	14903	0	12592	134	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 5.

All (134) 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:C:46[B]:ARG:NH2	9:C:401:HOH:O	2.05	0.90
3:G:138[B]:ASP:OD1	9:G:301:HOH:O	1.99	0.80
1:C:6[B]:ARG:NE	9:C:405:HOH:O	2.19	0.74
3:E:161:MET:HE3	3:E:166:PHE:CD1	2.23	0.74
4:F:67[A]:SER:OG	4:F:75[A]:SER:OG	2.05	0.74
4:F:68:ARG:NH2	9:F:301:HOH:O	2.00	0.74
1:C:230:ILE:HD11	4:H:206[B]:ARG:HD3	1.70	0.73
4:F:29:HIS:O	9:F:301:HOH:O	2.07	0.73
1:A:160:GLU:OE2	9:A:501:HOH:O	2.06	0.72
2:B:89:GLN:O	9:B:101:HOH:O	2.06	0.72
4:H:13:LEU:HD21	4:H:19[B]:MET:HG3	1.71	0.71
3:E:3[A]:ASN:OD1	9:E:301:HOH:O	2.08	0.70
1:C:155:ASN:OD1	3:E:66[B]:ARG:NH2	2.23	0.70
4:F:131:PRO:HD3	4:F:144:LEU:HG	1.73	0.69
1:C:58:HIS:ND1	9:C:407:HOH:O	2.26	0.68
4:F:65[A]:ASN:OD1	4:F:77:ARG:NH1	2.26	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:125:ASP:OD1	9:C:402:HOH:O	2.12	0.67
1:C:6[B]:ARG:HH11	1:C:8:PHE:HE1	1.40	0.67
4:F:13:LEU:HD21	4:F:19[B]:MET:HG3	1.76	0.67
3:E:161:MET:HE3	3:E:166:PHE:HD1	1.59	0.67
1:C:43:LYS:HD2	1:C:66[A]:LEU:CD1	2.26	0.66
3:E:3[A]:ASN:ND2	3:E:5:ASP:OD2	2.27	0.66
1:C:6[B]:ARG:NH1	9:C:410:HOH:O	2.29	0.65
3:G:3:ASN:ND2	3:G:5:ASP:OD2	2.24	0.65
1:C:43:LYS:HD3	1:C:62:TYR:HB3	1.79	0.65
4:F:111[A]:ARG:NH1	9:F:306:HOH:O	2.29	0.64
3:E:161:MET:CE	4:F:196:ARG:HD3	2.27	0.64
3:G:28:GLY:HA3	3:G:93[A]:SER:OG	1.99	0.62
3:G:80:MET:HG3	3:G:167:LYS:HE3	1.82	0.61
4:F:154[B]:ASP:OD1	9:F:302:HOH:O	2.16	0.61
4:F:5:THR:OG1	9:F:303:HOH:O	2.16	0.61
4:F:30:ASN:HD22	4:F:96:TRP:CG	2.19	0.61
3:E:66[B]:ARG:NH1	9:E:302:HOH:O	2.10	0.60
3:E:161:MET:HE2	4:F:196:ARG:HD3	1.83	0.60
4:F:71:LYS:NZ	9:F:307:HOH:O	2.30	0.59
3:E:25[A]:GLN:NE2	9:E:306:HOH:O	2.35	0.59
1:A:61[A]:ARG:HG2	3:G:94:ASN:HB3	1.84	0.59
1:C:188:ARG:NH2	9:C:406:HOH:O	2.21	0.59
1:C:43:LYS:HD2	1:C:66[A]:LEU:HD12	1.84	0.58
1:C:134:ASN:ND2	9:C:416:HOH:O	2.36	0.58
1:A:215:MET:HG3	1:A:257:HIS:CD2	2.38	0.58
3:E:3[A]:ASN:HD21	3:E:5:ASP:CG	2.12	0.58
1:A:77:LEU:HD13	1:A:92:TYR:HB2	1.86	0.57
1:C:246:LEU:O	9:C:403:HOH:O	2.17	0.57
3:E:176:ASN:O	3:E:176:ASN:ND2	2.37	0.57
3:E:157:CYS:SG	4:F:172[B]:CYS:HB2	2.45	0.56
1:A:113:ALA:HB2	2:B:60:TRP:CE2	2.40	0.55
3:E:28:GLY:HA3	3:E:93[A]:SER:OG	2.06	0.55
3:G:68:LYS:HD3	3:G:70[B]:TYR:HE2	1.71	0.55
1:A:64:GLN:OE1	9:A:502:HOH:O	2.19	0.54
4:F:28:ASN:HB2	9:F:422:HOH:O	2.07	0.54
4:F:71:LYS:HD3	9:F:359:HOH:O	2.07	0.53
1:C:222:VAL:HA	4:H:200[A]:THR:HG21	1.91	0.53
1:C:83:HIS:NE2	9:C:414:HOH:O	2.34	0.53
4:F:208:HIS:HB2	4:F:241:TRP:CZ3	2.44	0.52
4:F:116:GLU:OE1	9:F:304:HOH:O	2.19	0.52
3:E:145:GLN:NE2	9:E:310:HOH:O	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:111[A]:ARG:NH2	9:F:315:HOH:O	2.42	0.52
3:G:167:LYS:NZ	9:G:308:HOH:O	2.38	0.52
4:F:12:VAL:HG23	9:F:349:HOH:O	2.09	0.52
4:H:210[C]:ARG:HE	4:H:212:GLN:CG	2.23	0.52
1:C:211[B]:TYR:OH	9:C:404:HOH:O	2.19	0.51
4:H:212:GLN:NE2	9:H:403:HOH:O	2.27	0.51
1:A:134[A]:ASN:OD1	5:A:401:GOL:O2	2.29	0.51
4:F:194:ARG:N	4:F:194:ARG:HD2	2.26	0.51
4:F:172[B]:CYS:SG	4:F:194:ARG:HD3	2.50	0.51
1:A:43:LYS:HD3	1:A:62:TYR:HB3	1.92	0.50
1:C:113:ALA:HB2	2:D:60:TRP:CE2	2.46	0.50
3:E:55:GLU:HG3	9:H:404:HOH:O	2.12	0.49
4:H:8:PRO:HD2	4:H:21:LEU:HD22	1.94	0.49
1:C:221[B]:ILE:HG23	1:C:223:GLN:N	2.27	0.49
1:A:215:MET:HG3	1:A:257:HIS:NE2	2.27	0.49
2:D:21:ASN:HB3	2:D:70:PHE:CE1	2.47	0.49
1:C:1:ARG:HB2	9:C:535:HOH:O	2.12	0.49
1:A:72:MET:HG2	4:H:96:TRP:CD2	2.48	0.49
1:A:154:LYS:HD3	3:G:51:LEU:HD11	1.95	0.49
1:C:187:ASN:ND2	9:C:412:HOH:O	2.32	0.49
1:C:221[B]:ILE:HG23	1:C:223:GLN:H	1.78	0.48
4:F:19[A]:MET:HG2	4:F:78:LEU:HD13	1.95	0.48
3:E:50:VAL:O	3:E:66[A]:ARG:HD3	2.13	0.48
1:C:86:HIS:HB3	1:C:90:HIS:NE2	2.29	0.48
4:F:49[A]:SER:HB3	4:F:55:THR:HG22	1.96	0.48
1:A:34:THR:HB	1:A:43:LYS:HE3	1.96	0.48
4:F:26:ASP:OD1	4:F:26:ASP:N	2.43	0.47
4:F:11[A]:GLN:OE1	4:F:19[A]:MET:HE1	2.13	0.47
1:C:156:TRP:CZ3	1:C:161:CYS:HB2	2.50	0.47
1:C:77:LEU:HD13	1:C:92:TYR:HB2	1.96	0.47
5:D:102:GOL:O2	9:D:201:HOH:O	2.20	0.47
3:E:161:MET:HE1	4:F:196:ARG:HD3	1.96	0.47
3:G:166:PHE:CE2	4:H:141:LYS:HE2	2.51	0.46
4:H:19[B]:MET:HE1	4:H:21:LEU:HD21	1.96	0.46
4:H:210[C]:ARG:NH2	9:H:412:HOH:O	2.48	0.46
4:F:65[B]:ASN:ND2	9:F:319:HOH:O	2.46	0.46
3:G:28:GLY:HA3	3:G:93[B]:SER:HB3	1.98	0.46
1:C:123:ASN:ND2	9:C:431:HOH:O	2.48	0.46
4:H:128:VAL:HG23	4:H:238:ALA:HB3	1.96	0.46
4:F:3:GLY:HA3	4:F:27[A]:MET:HE3	1.97	0.45
4:F:154[A]:ASP:CG	4:F:177:PRO:HG3	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:19[B]:MET:HE3	4:F:19[B]:MET:HB3	1.89	0.45
1:A:43:LYS:NZ	9:A:510:HOH:O	2.46	0.45
4:F:188:ARG:NH2	9:F:324:HOH:O	2.50	0.45
3:E:28:GLY:HA3	3:E:93[B]:SER:HB3	1.99	0.45
1:A:43:LYS:HD2	1:A:66[A]:LEU:HD12	1.99	0.44
2:B:72:PRO:HB2	2:B:97:ARG:NH2	2.31	0.44
3:E:3[A]:ASN:HB3	3:E:25[A]:GLN:HG3	1.99	0.44
4:F:30:ASN:O	4:F:51:SER:HA	2.17	0.44
4:H:210[C]:ARG:NH2	4:H:212:GLN:OE1	2.50	0.44
3:E:36:GLN:HB2	3:E:46:LEU:HD11	1.99	0.44
3:E:152:TYR:O	3:E:173:ALA:HA	2.16	0.44
1:C:43:LYS:HD2	1:C:66[A]:LEU:HD11	1.98	0.44
1:C:186:VAL:HG11	1:C:269:VAL:HG22	1.99	0.44
1:C:36:TYR:CB	1:C:66[A]:LEU:HD13	2.47	0.44
4:F:25:GLN:O	4:F:72[B]:ARG:HG2	2.18	0.43
3:G:166:PHE:CD2	4:H:141:LYS:HE2	2.53	0.43
1:C:72:MET:HE2	9:C:656:HOH:O	2.18	0.43
4:F:6:GLN:HA	4:F:22:GLN:O	2.18	0.43
4:H:3:GLY:HA2	4:H:27[A]:MET:HE3	2.00	0.43
3:E:159:LEU:HD11	3:E:161:MET:HE2	2.01	0.42
4:H:224:TRP:CZ2	4:H:226:GLN:HB2	2.54	0.42
4:H:19[B]:MET:HE1	4:H:112:LEU:HD11	2.01	0.42
4:F:21:LEU:HB2	4:F:76:LEU:HB3	2.02	0.41
4:F:30:ASN:HB2	4:F:95:VAL:O	2.19	0.41
4:F:155:HIS:HB3	4:F:216:TYR:HB2	2.02	0.41
3:E:117:ALA:HB2	3:E:196:PHE:HB3	2.02	0.41
4:F:94:SER:HB3	4:F:104:LEU:HD23	2.01	0.41
1:A:3:HIS:CG	1:A:169:LEU:HD21	2.56	0.41
1:A:43:LYS:HD2	1:A:66[A]:LEU:CD1	2.50	0.41
1:C:169[A]:LEU:HD23	1:C:176:LEU:HD13	2.02	0.41
4:F:212:GLN:HG3	4:F:235:ILE:HG23	2.03	0.41
1:C:198:LEU:HD11	1:C:246:LEU:HD12	2.02	0.40
4:F:47:TYR:HB2	4:F:66:VAL:HG11	2.02	0.40
1:A:36:TYR:CB	1:A:66[A]:LEU:HD13	2.51	0.40
4:H:117[B]:ASP:OD1	4:H:119:LYS:HG3	2.21	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)
9:C:448:HOH:O	9:F:449:HOH:O[4_443]	2.16	0.04

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	271/271 (100%)	264 (97%)	7 (3%)	0	100	100
1	C	280/271 (103%)	275 (98%)	5 (2%)	0	100	100
2	B	97/100 (97%)	96 (99%)	1 (1%)	0	100	100
2	D	101/100 (101%)	100 (99%)	1 (1%)	0	100	100
3	E	198/204 (97%)	193 (98%)	5 (2%)	0	100	100
3	G	213/204 (104%)	208 (98%)	5 (2%)	0	100	100
4	F	261/246 (106%)	256 (98%)	5 (2%)	0	100	100
4	H	254/246 (103%)	248 (98%)	6 (2%)	0	100	100
All	All	1675/1642 (102%)	1640 (98%)	35 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	233/241 (97%)	232 (100%)	1 (0%)	84	85
1	C	249/241 (103%)	244 (98%)	5 (2%)	48	43
2	B	82/95 (86%)	81 (99%)	1 (1%)	63	61
2	D	92/95 (97%)	89 (97%)	3 (3%)	33	24
3	E	171/181 (94%)	166 (97%)	5 (3%)	37	29
3	G	188/181 (104%)	184 (98%)	4 (2%)	47	41

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	F	219/212 (103%)	213 (97%)	6 (3%)	39	32
4	H	213/212 (100%)	203 (95%)	10 (5%)	23	12
All	All	1447/1458 (99%)	1412 (98%)	35 (2%)	48	36

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	MET
2	B	97	ARG
1	C	16	ILE
1	C	17	HIS
1	C	80	LEU
1	C	188	ARG
1	C	259	GLU
2	D	64[A]	LEU
2	D	64[B]	LEU
2	D	99	MET
3	E	25[A]	GLN
3	E	25[B]	GLN
3	E	157	CYS
3	E	188	ASN
3	E	194	ASP
4	F	11[A]	GLN
4	F	11[C]	GLN
4	F	19[A]	MET
4	F	19[B]	MET
4	F	185	ASN
4	F	194	ARG
3	G	4	ILE
3	G	112	GLN
3	G	146[A]	SER
3	G	146[B]	SER
4	H	18	SER
4	H	65[A]	ASN
4	H	65[B]	ASN
4	H	67	SER
4	H	120[A]	ASN
4	H	120[B]	ASN
4	H	132	SER
4	H	172[A]	CYS
4	H	172[B]	CYS

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Mol	Chain	Res	Type
4	H	184	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	137	HIS
1	C	187	ASN
1	C	203	HIS
2	D	13	HIS
2	D	17	ASN
3	E	19	GLN
3	E	36	GLN
3	E	188	ASN
3	G	49	ASN
3	G	79	GLN
3	G	94	ASN
3	G	120	GLN
4	H	11	GLN
4	H	214	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 1 is 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
5	GOL	D	102	-	5,5,5	1.24	1 (20%)	5,5,5	0.97	0
6	P1J	A	402	-	16,16,16	1.07	2 (12%)	20,21,21	1.10	2 (10%)
6	P1J	C	301	-	16,16,16	1.18	2 (12%)	20,21,21	1.01	2 (10%)
5	GOL	A	401	-	5,5,5	1.29	1 (20%)	5,5,5	1.54	1 (20%)
8	ACT	H	301	-	3,3,3	1.64	1 (33%)	3,3,3	1.34	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
6	P1J	C	301	-	-	2/10/10/10	0/1/1/1
5	GOL	D	102	-	-	0/4/4/4	-
6	P1J	A	402	-	-	3/10/10/10	0/1/1/1
5	GOL	A	401	-	-	2/4/4/4	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	301	P1J	O-C	-3.20	1.20	1.30
6	C	301	P1J	O1-C	3.16	1.32	1.22
6	A	402	P1J	O1-C	2.96	1.31	1.22
6	A	402	P1J	O-C	-2.89	1.21	1.30
8	H	301	ACT	CH3-C	2.42	1.58	1.49
5	A	401	GOL	O2-C2	-2.37	1.36	1.43
5	D	102	GOL	C1-C2	2.01	1.59	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	402	P1J	O1-C-C1	-3.32	112.57	123.09
6	A	402	P1J	O-C-C1	3.07	123.71	114.00
6	C	301	P1J	O1-C-C1	-2.58	114.92	123.09
6	C	301	P1J	O-C-C1	2.18	120.87	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	401	GOL	C3-C2-C1	-2.08	104.17	111.80

There are no chirality outliers.

All (7) torsion outliers are listed below:

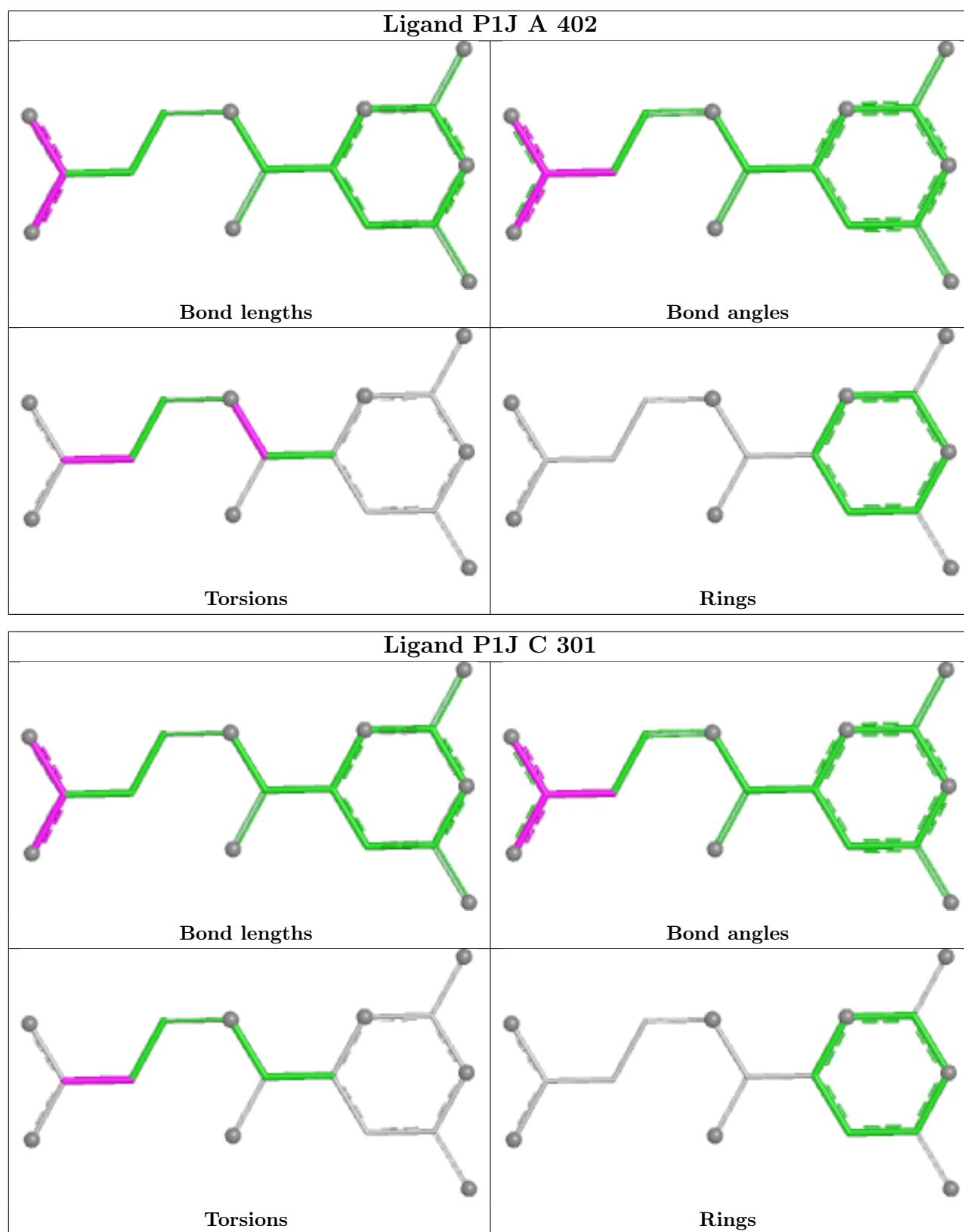
Mol	Chain	Res	Type	Atoms
5	A	401	GOL	O1-C1-C2-C3
5	A	401	GOL	O1-C1-C2-O2
6	C	301	P1J	O-C-C1-C2
6	C	301	P1J	O1-C-C1-C2
6	A	402	P1J	O-C-C1-C2
6	A	402	P1J	O2-C3-N-C2
6	A	402	P1J	O1-C-C1-C2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	102	GOL	1	0
5	A	401	GOL	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 ⓘ

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	265/271 (97%)	0.12	6 (2%) 61 68	18, 42, 81, 122	16 (6%)
1	C	267/271 (98%)	-0.10	8 (2%) 52 59	14, 34, 61, 97	27 (10%)
2	B	98/100 (98%)	0.63	4 (4%) 41 48	31, 63, 91, 98	2 (2%)
2	D	100/100 (100%)	-0.12	1 (1%) 79 84	20, 40, 62, 68	10 (10%)
3	E	192/204 (94%)	0.07	1 (0%) 87 90	15, 42, 86, 102	16 (8%)
3	G	201/204 (98%)	-0.15	3 (1%) 72 78	14, 35, 58, 110	21 (10%)
4	F	242/246 (98%)	0.58	26 (10%) 11 12	19, 47, 74, 99	34 (14%)
4	H	242/246 (98%)	0.04	5 (2%) 63 70	15, 41, 66, 88	20 (8%)
All	All	1607/1642 (97%)	0.12	54 (3%) 48 54	14, 41, 77, 122	146 (9%)

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	F	50	ALA	4.8
1	C	221[A]	ILE	4.4
1	A	222	VAL	4.3
4	F	49[A]	SER	4.1
4	F	244	ALA	4.0
4	F	96	TRP	3.9
4	F	98	GLY	3.8
4	F	69	LEU	3.7
4	F	70	ASN	3.2
1	C	16	ILE	3.2
1	A	221	ILE	3.1
4	F	97	THR	3.1
4	F	52[A]	GLU	3.1
2	B	49	VAL	3.0
1	C	219	GLU	2.7
2	D	99	MET	2.7

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Mol	Chain	Res	Type	RSRZ
3	E	130	SER	2.7
1	A	249	GLN	2.7
1	C	222	VAL	2.7
4	F	74	PHE	2.7
1	A	61[A]	ARG	2.7
1	A	246	LEU	2.6
4	H	53	GLY	2.5
4	F	67[A]	SER	2.5
4	F	54	THR	2.5
1	C	247	ASP	2.4
4	F	30	ASN	2.4
4	H	65[A]	ASN	2.4
4	F	3	GLY	2.4
4	H	47	TYR	2.3
1	C	253	LEU	2.3
4	H	244	ALA	2.3
3	G	200	PRO	2.3
3	G	202	SER	2.3
4	F	53	GLY	2.3
1	C	246	LEU	2.3
4	F	241	TRP	2.2
4	F	51	SER	2.2
2	B	0	MET	2.2
4	F	72[A]	ARG	2.2
1	C	211[A]	TYR	2.1
4	F	55	THR	2.1
4	F	31	SER	2.1
2	B	51	HIS	2.1
4	F	27[A]	MET	2.1
1	A	55	ALA	2.1
4	F	26	ASP	2.1
4	H	26	ASP	2.1
3	G	201	GLU	2.1
4	F	73	GLU	2.1
4	F	28	ASN	2.1
4	F	221	ASN	2.1
2	B	89	GLN	2.1
4	F	77	ARG	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

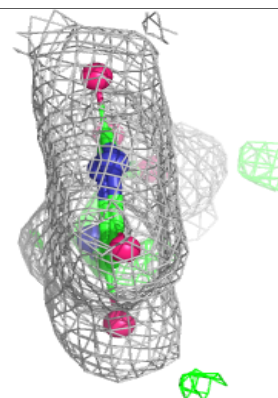
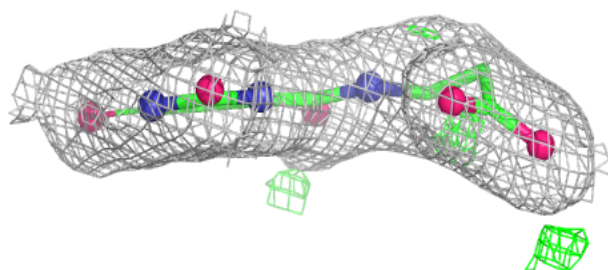
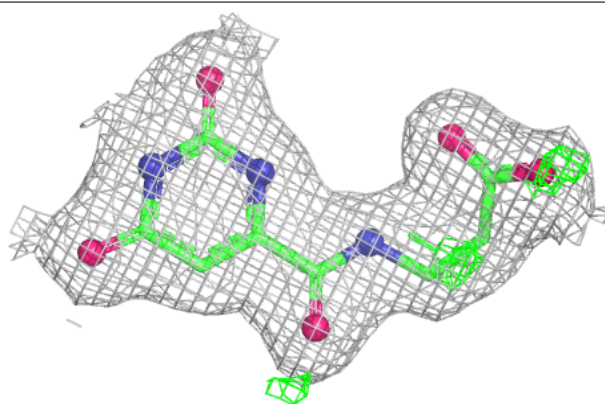
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	ACT	H	301	4/4	0.84	0.14	66,68,69,69	0
5	GOL	A	401	6/6	0.94	0.11	38,50,53,55	0
5	GOL	D	102	6/6	0.95	0.11	36,38,40,42	0
6	P1J	C	301	16/16	0.97	0.05	24,29,37,39	0
7	NA	D	101	1/1	0.97	0.11	46,46,46,46	0
6	P1J	A	402	16/16	0.97	0.05	27,34,42,42	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.

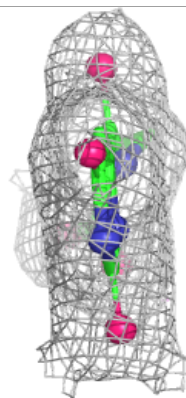
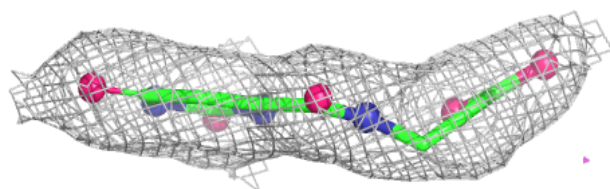
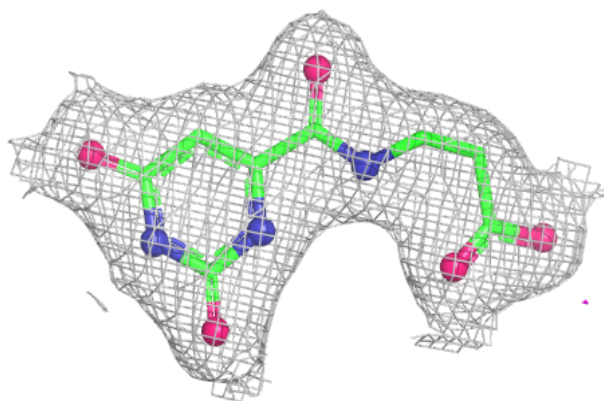
Electron density around P1J C 301:

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 P1J A 402:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.