



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:22 PM UTC

PDB ID : 6B3E / pdb_00006b3e
Title : Crystal structure of human CDK12/CyclinK in complex with an inhibitor
Authors : Ferguson, A.D.
Deposited on : 2017-09-21
Resolution : 3.06 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

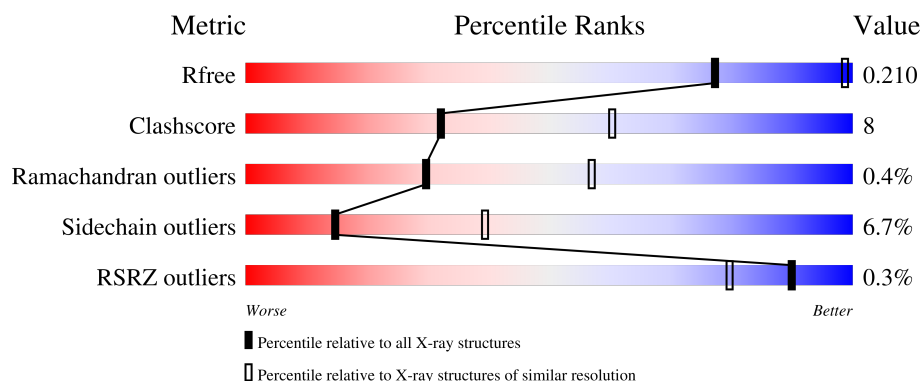
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.06 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	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	320	 73% 20% . .
1	C	320	 72% 21% . .
2	B	267	 78% 10% . 10%
2	D	267	 78% 10% . 10%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9119 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 Cyclin-dependent kinase 12.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	310	Total	C	N	O	P	S	0	0	0
			2521	1621	425	458	1	16			
1	C	308	Total	C	N	O	P	S	0	0	0
			2507	1614	423	453	1	16			

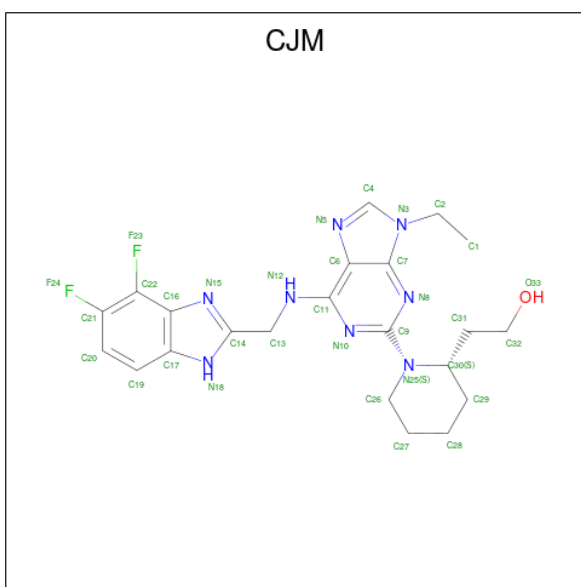
- Molecule 2 is a protein called Cyclin-K.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	240	Total	C	N	O	S	0	0	0
			1998	1302	332	352	12			
2	D	241	Total	C	N	O	S	0	0	0
			2004	1305	334	353	12			

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

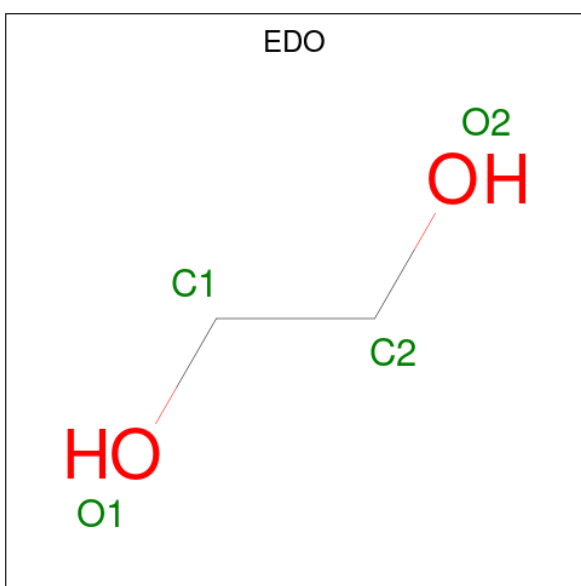
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is 2-[(2S)-1-(6-[[[(4,5-difluoro-1H-benzimidazol-2-yl)methyl]amino}-9-ethyl-9H-purin-2-yl]piperidin-2-yl]ethan-1-ol (CCD ID: CJM) (formula: C₂₂H₂₆F₂N₈O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	F	N	O	0	0
			33	22	2	8	1		
4	C	1	Total	C	F	N	O	0	0
			33	22	2	8	1		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		

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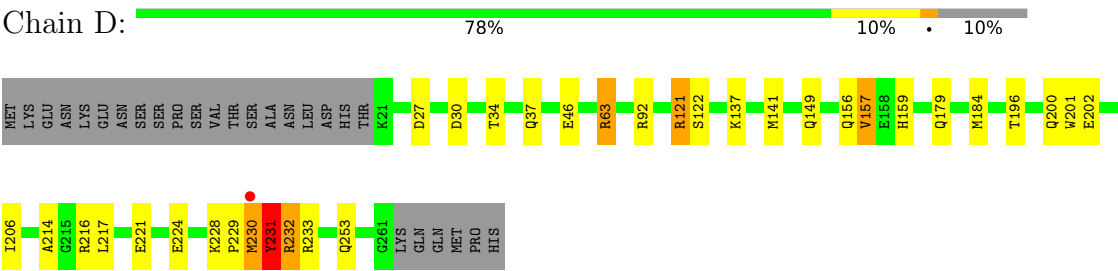
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	6	Total	O	0	0
			6	6		
6	B	1	Total	O	0	0
			1	1		
6	D	2	Total	O	0	0
			2	2		

● Molecule 2: Cyclin-K



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	50.14Å 77.26Å 90.64Å 76.53° 85.28° 78.22°	Depositor
Resolution (Å)	88.08 – 3.06 88.08 – 3.06	Depositor EDS
% Data completeness (in resolution range)	93.4 (88.08-3.06) 93.4 (88.08-3.06)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 3.07Å)	Xtriage
Refinement program	BUSTER	Depositor
R, R_{free}	0.190 , 0.212 0.190 , 0.210	Depositor DCC
R_{free} test set	1111 reflections (4.52%)	wwPDB-VP
Wilson B-factor (Å ²)	73.0	Xtriage
Anisotropy	0.646	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 86.3	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	0.95	EDS
Total number of atoms	9119	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.26% 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: TPO, CJM, MG, EDO

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	0.79	1/2566 (0.0%)	1.13	3/3459 (0.1%)
1	C	0.76	2/2552 (0.1%)	1.16	6/3440 (0.2%)
2	B	0.77	0/2053	1.12	3/2779 (0.1%)
2	D	0.75	0/2059	1.21	8/2786 (0.3%)
All	All	0.77	3/9230 (0.0%)	1.16	20/12464 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	821	MET	SD-CE	6.17	1.95	1.79
1	A	821	MET	SD-CE	5.90	1.94	1.79
1	C	844	MET	SD-CE	-5.38	1.66	1.79

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	232	ARG	CA-C-N	10.67	141.91	121.54
2	D	232	ARG	C-N-CA	10.67	141.91	121.54
2	D	230	MET	CA-C-N	10.61	141.80	121.54
2	D	230	MET	C-N-CA	10.61	141.80	121.54
1	C	760	LEU	N-CA-C	-10.20	99.94	114.12
2	D	231	TYR	CA-C-N	7.34	135.56	121.54
2	D	231	TYR	C-N-CA	7.34	135.56	121.54
1	C	763	GLU	CA-C-N	6.77	131.84	120.88
1	C	763	GLU	C-N-CA	6.77	131.84	120.88
1	C	1035	HIS	CA-C-N	6.64	133.65	121.70
1	C	1035	HIS	C-N-CA	6.64	133.65	121.70
1	A	1022	ASP	CA-CB-CG	5.91	118.51	112.60
1	C	725	ASP	CA-CB-CG	5.77	118.37	112.60
1	A	1035	HIS	N-CA-C	-5.75	106.17	112.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	27	ASP	CA-CB-CG	5.74	118.34	112.60
1	A	725	ASP	CA-CB-CG	5.70	118.30	112.60
2	B	126	ASP	CA-CB-CG	5.56	118.16	112.60
2	B	92	ARG	N-CA-C	5.31	117.06	111.28
2	D	92	ARG	N-CA-C	5.13	116.87	111.28
2	D	231	TYR	N-CA-CB	-5.07	101.91	110.49

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	2521	0	2548	53	0
1	C	2507	0	2539	59	0
2	B	1998	0	1985	21	0
2	D	2004	0	1989	22	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
4	A	33	0	0	4	0
4	C	33	0	0	0	0
5	B	8	0	12	1	0
5	C	4	0	6	1	0
6	A	6	0	0	0	0
6	B	1	0	0	0	0
6	D	2	0	0	0	0
All	All	9119	0	9079	146	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 8.

All (146) 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:739:GLY:HA2	1:C:759:ARG:NH1	1.57	1.17

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:999:HIS:CE1	1:A:1009:CYS:HB3	2.06	0.90
1:A:733:ILE:HG22	4:A:1102:CJM:C20	2.02	0.90
1:C:999:HIS:CE1	1:C:1009:CYS:HB3	2.08	0.89
1:A:960:TRP:O	1:A:963:VAL:HG12	1.76	0.86
1:C:1024:GLU:OE1	1:C:1027:LYS:HD2	1.80	0.81
1:A:796:LYS:H	1:A:796:LYS:HD3	1.44	0.81
1:C:960:TRP:O	1:C:963:VAL:HG12	1.80	0.80
1:A:796:LYS:HD3	1:A:796:LYS:N	1.97	0.79
1:A:757:LYS:HD3	1:A:759:ARG:HE	1.48	0.78
1:C:796:LYS:HD3	1:C:796:LYS:H	1.49	0.77
1:A:1024:GLU:OE1	1:A:1027:LYS:HD2	1.85	0.77
1:C:739:GLY:HA2	1:C:759:ARG:HH11	1.52	0.72
2:B:179:GLN:HE22	1:C:1007:LYS:NZ	1.87	0.72
1:A:733:ILE:HG22	4:A:1102:CJM:C19	2.21	0.71
1:C:796:LYS:HD3	1:C:796:LYS:N	2.06	0.70
1:A:722:ARG:HE	1:A:726:LYS:HD3	1.56	0.69
1:A:834:ASP:O	1:A:1028:MET:HE2	1.94	0.68
1:C:722:ARG:HE	1:C:726:LYS:HD3	1.58	0.68
1:A:986:PHE:HB3	1:A:989:ILE:HD12	1.76	0.66
2:B:175:LYS:HZ3	1:C:999:HIS:CE1	2.12	0.66
1:A:722:ARG:HG2	1:A:726:LYS:NZ	2.10	0.66
1:C:986:PHE:HB3	1:C:989:ILE:HD12	1.76	0.66
2:D:121:ARG:O	2:D:121:ARG:HD3	1.95	0.66
2:B:184:MET:HE1	2:B:217:LEU:HB2	1.78	0.66
2:D:184:MET:HE1	2:D:217:LEU:HB2	1.78	0.66
1:C:722:ARG:HG2	1:C:726:LYS:NZ	2.11	0.65
1:A:722:ARG:HE	1:A:726:LYS:CD	2.10	0.64
2:B:175:LYS:NZ	1:C:999:HIS:CE1	2.66	0.63
1:A:733:ILE:CG2	4:A:1102:CJM:C19	2.78	0.62
1:C:722:ARG:HE	1:C:726:LYS:CD	2.13	0.62
2:B:184:MET:HE2	2:B:214:ALA:HA	1.82	0.62
2:D:27:ASP:HB3	2:D:30:ASP:H	1.65	0.61
1:C:719:TRP:HZ3	1:C:789:MET:HE3	1.66	0.60
1:C:739:GLY:HA2	1:C:759:ARG:HH12	1.55	0.60
1:A:719:TRP:HZ3	1:A:789:MET:HE3	1.67	0.59
1:C:960:TRP:O	1:C:963:VAL:CG1	2.49	0.59
1:C:832:SER:H	1:C:835:HIS:CD2	2.21	0.59
1:A:796:LYS:H	1:A:796:LYS:CD	2.16	0.58
1:C:840:MET:HE3	1:C:844:MET:HE3	1.85	0.58
1:A:769:ILE:O	1:A:773:ARG:HG3	2.03	0.58
1:A:840:MET:HE3	1:A:844:MET:HE3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:184:MET:HE2	2:D:214:ALA:HA	1.85	0.57
1:A:976:LYS:HG2	1:A:977:GLN:N	2.20	0.57
1:C:769:ILE:O	1:C:773:ARG:HG3	2.05	0.57
1:A:832:SER:H	1:A:835:HIS:CD2	2.23	0.57
1:A:840:MET:CE	1:A:844:MET:HE3	2.35	0.56
2:B:20:THR:HG23	2:B:21:LYS:H	1.70	0.56
2:B:200:GLN:HE22	2:B:253:GLN:HE22	1.54	0.56
2:D:34:THR:OG1	2:D:37:GLN:HG2	2.05	0.56
1:C:999:HIS:ND1	1:C:1009:CYS:HB3	2.21	0.56
1:C:840:MET:CE	1:C:844:MET:HE3	2.36	0.55
2:B:63:ARG:NH2	2:B:122:SER:OG	2.38	0.55
1:C:774:GLU:HG2	1:C:778:LEU:HD12	1.87	0.55
2:D:200:GLN:HE22	2:D:253:GLN:HE22	1.52	0.55
1:A:722:ARG:HH21	1:A:726:LYS:HB3	1.72	0.55
1:A:999:HIS:ND1	1:A:1009:CYS:HB3	2.22	0.54
1:A:854:ASN:O	1:A:883:LEU:HD22	2.08	0.54
1:C:760:LEU:O	1:C:763:GLU:HB2	2.08	0.54
2:D:121:ARG:HD3	2:D:121:ARG:C	2.32	0.54
2:D:184:MET:CE	2:D:214:ALA:HA	2.38	0.54
2:B:179:GLN:HE22	1:C:1007:LYS:HZ2	1.52	0.54
2:B:184:MET:CE	2:B:214:ALA:HA	2.38	0.54
1:C:931:THR:HG22	1:C:988:PHE:HZ	1.72	0.54
1:A:1007:LYS:HZ2	2:D:179:GLN:HE22	1.56	0.54
2:B:69:ASP:H	5:B:301:EDO:H12	1.72	0.54
1:C:877:ASP:H	5:C:1102:EDO:H22	1.73	0.53
1:A:775:ILE:O	1:A:779:ARG:HG3	2.08	0.53
1:C:722:ARG:HH21	1:C:726:LYS:HB3	1.72	0.53
2:D:63:ARG:NH2	2:D:122:SER:OG	2.42	0.53
2:D:27:ASP:HB2	2:D:30:ASP:OD2	2.10	0.52
1:A:931:THR:HG22	1:A:988:PHE:HZ	1.73	0.52
1:C:722:ARG:HG2	1:C:726:LYS:HZ2	1.75	0.52
1:A:1007:LYS:NZ	2:D:179:GLN:HE22	2.07	0.52
1:C:970:ASN:HA	1:C:973:LYS:HE3	1.92	0.51
1:C:936:PHE:CB	1:C:944:GLN:HE22	2.23	0.51
1:C:796:LYS:H	1:C:796:LYS:CD	2.21	0.51
1:A:936:PHE:CB	1:A:944:GLN:HE22	2.24	0.51
1:A:970:ASN:HA	1:A:973:LYS:HE3	1.93	0.50
1:A:936:PHE:HB2	1:A:944:GLN:HE22	1.76	0.50
1:A:842:GLN:NE2	1:A:873:ILE:H	2.08	0.50
1:C:842:GLN:NE2	1:C:873:ILE:H	2.10	0.50
1:C:936:PHE:HB2	1:C:944:GLN:HE22	1.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:760:LEU:HG	1:A:807:GLY:O	2.12	0.49
2:B:137:LYS:O	2:B:141:MET:HG2	2.13	0.49
1:C:821:MET:HE3	1:C:863:SER:HB3	1.93	0.49
1:A:844:MET:HE1	1:A:922:CYS:HB3	1.94	0.49
2:B:34:THR:OG1	2:B:37:GLN:HG2	2.12	0.49
1:A:722:ARG:HG2	1:A:726:LYS:HZ3	1.78	0.48
1:C:1021:LYS:CD	1:C:1022:ASP:CG	2.86	0.48
1:C:775:ILE:O	1:C:779:ARG:HG3	2.12	0.48
1:A:719:TRP:CZ3	1:A:789:MET:HE3	2.49	0.48
2:D:137:LYS:O	2:D:141:MET:HG2	2.13	0.48
1:A:815:TYR:OH	4:A:1102:CJM:N18	2.46	0.48
2:B:184:MET:CE	2:B:217:LEU:HB2	2.44	0.48
1:A:936:PHE:HB3	1:A:944:GLN:NE2	2.29	0.47
1:C:936:PHE:HB3	1:C:944:GLN:NE2	2.29	0.47
2:D:184:MET:CE	2:D:217:LEU:HB2	2.43	0.47
1:A:821:MET:HE3	1:A:863:SER:HB3	1.97	0.47
1:A:889:SER:O	1:A:890:ARG:HD3	2.14	0.47
1:A:722:ARG:HG2	1:A:726:LYS:HZ2	1.78	0.47
1:C:722:ARG:NH2	1:C:726:LYS:HB3	2.29	0.47
1:C:844:MET:HE1	1:C:922:CYS:HB3	1.97	0.47
1:A:782:ILE:HD12	1:A:849:TYR:OH	2.13	0.47
1:C:841:LYS:HG2	1:C:1025:LEU:HD11	1.97	0.47
1:A:976:LYS:CG	1:A:977:GLN:N	2.78	0.46
2:B:157:VAL:HG12	2:B:159:HIS:CE1	2.51	0.46
2:D:157:VAL:HG12	2:D:159:HIS:CE1	2.50	0.46
2:B:184:MET:HE2	2:B:214:ALA:CA	2.45	0.46
2:B:201:TRP:HB2	2:B:206:ILE:HD11	1.98	0.46
1:A:842:GLN:HE22	1:A:873:ILE:H	1.63	0.45
1:C:721:LYS:HD3	2:D:149:GLN:HB3	1.98	0.45
1:C:719:TRP:CZ3	1:C:789:MET:HE3	2.49	0.45
2:D:201:TRP:HB2	2:D:206:ILE:HD11	1.97	0.45
1:C:914:THR:HB	1:C:915:PRO:HD2	1.98	0.45
1:A:722:ARG:NH2	1:A:726:LYS:HB3	2.31	0.45
1:C:741:VAL:HG22	1:C:756:LYS:HG3	1.98	0.45
1:A:841:LYS:HE3	1:A:845:GLU:OE2	2.17	0.44
1:A:834:ASP:O	1:A:1028:MET:CE	2.64	0.44
1:C:745:LYS:HE2	1:C:752:LEU:HD13	2.00	0.44
1:A:914:THR:HB	1:A:915:PRO:HD2	2.00	0.44
2:D:184:MET:HE2	2:D:214:ALA:CA	2.47	0.44
1:A:936:PHE:CB	1:A:944:GLN:NE2	2.81	0.44
1:C:976:LYS:CG	1:C:977:GLN:N	2.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:945:LEU:HD22	1:C:972:MET:HE2	1.99	0.43
1:C:936:PHE:CB	1:C:944:GLN:NE2	2.81	0.43
1:C:842:GLN:HE22	1:C:873:ILE:H	1.65	0.43
1:A:914:THR:HB	1:A:915:PRO:CD	2.49	0.42
1:C:841:LYS:HG2	1:C:1025:LEU:HD21	2.01	0.42
1:C:960:TRP:CG	1:C:963:VAL:HB	2.54	0.42
1:A:756:LYS:HG2	1:A:811:LEU:HD22	2.00	0.42
1:C:914:THR:HB	1:C:915:PRO:CD	2.49	0.42
2:D:196:THR:O	2:D:200:GLN:HG3	2.20	0.42
1:A:1035:HIS:CD2	1:A:1035:HIS:N	2.87	0.42
1:C:889:SER:O	1:C:890:ARG:HD3	2.20	0.42
2:D:233:ARG:HD2	2:D:233:ARG:HA	1.79	0.41
2:B:179:GLN:HE22	1:C:1007:LYS:HZ3	1.62	0.41
2:D:229:PRO:HB2	2:D:231:TYR:HB3	2.03	0.41
1:C:821:MET:CE	1:C:863:SER:HB3	2.50	0.41
1:A:946:GLU:O	1:A:950:ARG:HG3	2.20	0.41
2:B:37:GLN:HG2	2:B:37:GLN:H	1.73	0.41
2:B:179:GLN:NE2	1:C:1007:LYS:NZ	2.62	0.41
1:C:722:ARG:HG2	1:C:726:LYS:HZ3	1.84	0.41
1:C:841:LYS:HE3	1:C:845:GLU:OE2	2.21	0.41
2:B:223:GLN:O	2:B:229:PRO:HA	2.21	0.40
2:D:37:GLN:NE2	2:D:46:GLU:OE1	2.48	0.40

There are no symmetry-related clashes.

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	A	305/320 (95%)	299 (98%)	5 (2%)	1 (0%)	36	63
1	C	303/320 (95%)	295 (97%)	7 (2%)	1 (0%)	36	63
2	B	238/267 (89%)	237 (100%)	1 (0%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	239/267 (90%)	235 (98%)	2 (1%)	2 (1%)	16	42
All	All	1085/1174 (92%)	1066 (98%)	15 (1%)	4 (0%)	30	57

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	231	TYR
2	D	232	ARG
1	A	897	ILE
1	C	897	ILE

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	278/287 (97%)	255 (92%)	23 (8%)	10	32
1	C	276/287 (96%)	253 (92%)	23 (8%)	10	32
2	B	216/241 (90%)	206 (95%)	10 (5%)	24	52
2	D	216/241 (90%)	206 (95%)	10 (5%)	24	52
All	All	986/1056 (93%)	920 (93%)	66 (7%)	15	39

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	717	SER
1	A	722	ARG
1	A	735	GLU
1	A	752	LEU
1	A	756	LYS
1	A	760	LEU
1	A	782	ILE
1	A	784	ARG
1	A	787	VAL
1	A	796	LYS

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Mol	Chain	Res	Type
1	A	820	LEU
1	A	828	LEU
1	A	829	VAL
1	A	908	LEU
1	A	942	LEU
1	A	977	GLN
1	A	979	ARG
1	A	984	GLU
1	A	988	PHE
1	A	1020	LEU
1	A	1022	ASP
1	A	1025	LEU
1	A	1028	MET
2	B	37	GLN
2	B	63	ARG
2	B	126	ASP
2	B	156	GLN
2	B	157	VAL
2	B	202	GLU
2	B	216	ARG
2	B	221	GLU
2	B	224	GLU
2	B	232	ARG
1	C	735	GLU
1	C	752	LEU
1	C	760	LEU
1	C	764	LYS
1	C	776	LYS
1	C	784	ARG
1	C	787	VAL
1	C	796	LYS
1	C	828	LEU
1	C	829	VAL
1	C	908	LEU
1	C	929	LEU
1	C	942	LEU
1	C	962	ASP
1	C	977	GLN
1	C	979	ARG
1	C	984	GLU
1	C	988	PHE
1	C	1020	LEU

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Mol	Chain	Res	Type
1	C	1021	LYS
1	C	1022	ASP
1	C	1028	MET
1	C	1036	TRP
2	D	63	ARG
2	D	121	ARG
2	D	156	GLN
2	D	157	VAL
2	D	202	GLU
2	D	216	ARG
2	D	221	GLU
2	D	224	GLU
2	D	228	LYS
2	D	230	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	780	GLN
1	A	835	HIS
1	A	842	GLN
1	A	872	GLN
1	A	944	GLN
1	A	999	HIS
1	A	1035	HIS
2	B	149	GLN
2	B	179	GLN
2	B	200	GLN
1	C	835	HIS
1	C	842	GLN
1	C	944	GLN
1	C	999	HIS
2	D	33	HIS
2	D	179	GLN
2	D	200	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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
1	TPO	A	893	1	8,10,11	1.33	1 (12%)	10,14,16	1.92	3 (30%)
1	TPO	C	893	1	8,10,11	1.03	0	10,14,16	1.78	2 (20%)

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
1	TPO	A	893	1	-	1/9/11/13	-
1	TPO	C	893	1	-	0/9/11/13	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	893	TPO	CB-CA	2.74	1.59	1.53

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	893	TPO	CG2-CB-CA	-4.15	105.17	113.26
1	C	893	TPO	CG2-CB-CA	-3.92	105.63	113.26
1	A	893	TPO	O-C-CA	-2.28	118.91	124.77
1	A	893	TPO	O3P-P-OG1	2.08	113.94	105.85
1	C	893	TPO	O-C-CA	-2.05	119.51	124.77

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	893	TPO	O-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	CJM	A	1102	-	37,37,37	1.22	2 (5%)	47,53,53	1.59	7 (14%)
4	CJM	C	1103	-	37,37,37	1.31	3 (8%)	47,53,53	1.69	5 (10%)
5	EDO	B	301	-	3,3,3	0.37	0	2,2,2	0.56	0
5	EDO	C	1102	-	3,3,3	0.71	0	2,2,2	0.24	0
5	EDO	B	302	-	3,3,3	0.77	0	2,2,2	0.14	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
4	CJM	A	1102	-	-	4/14/25/25	1/5/5/5
4	CJM	C	1103	-	-	0/14/25/25	1/5/5/5
5	EDO	B	301	-	-	1/1/1/1	-
5	EDO	C	1102	-	-	0/1/1/1	-
5	EDO	B	302	-	-	0/1/1/1	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1103	CJM	C13-C14	5.79	1.53	1.49
4	A	1102	CJM	C13-C14	-4.88	1.46	1.49
4	C	1103	CJM	C9-N25	3.97	1.43	1.36
4	A	1102	CJM	C17-C16	-3.86	1.36	1.40
4	C	1103	CJM	C14-N15	2.49	1.36	1.32

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1102	CJM	C16-C17-N18	7.61	109.46	105.13
4	C	1103	CJM	C16-C17-N18	7.53	109.42	105.13
4	C	1103	CJM	C17-C16-N15	-5.67	105.40	110.31
4	A	1102	CJM	C17-C16-N15	-4.40	106.50	110.31
4	C	1103	CJM	C19-C17-C16	-3.04	119.71	122.57
4	A	1102	CJM	C13-C14-N18	2.93	126.82	122.72
4	A	1102	CJM	C13-C14-N15	-2.80	120.76	124.41
4	C	1103	CJM	C22-C16-N15	2.67	135.74	129.03
4	C	1103	CJM	C26-N25-C30	2.49	118.48	115.20
4	A	1102	CJM	C19-C17-C16	-2.48	120.24	122.57
4	A	1102	CJM	C17-N18-C14	-2.16	104.73	107.23
4	A	1102	CJM	C22-C16-N15	2.06	134.20	129.03

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1102	CJM	N25-C30-C31-C32
4	A	1102	CJM	C29-C30-C31-C32
5	B	301	EDO	O1-C1-C2-O2
4	A	1102	CJM	N12-C13-C14-N18
4	A	1102	CJM	C30-C31-C32-O33

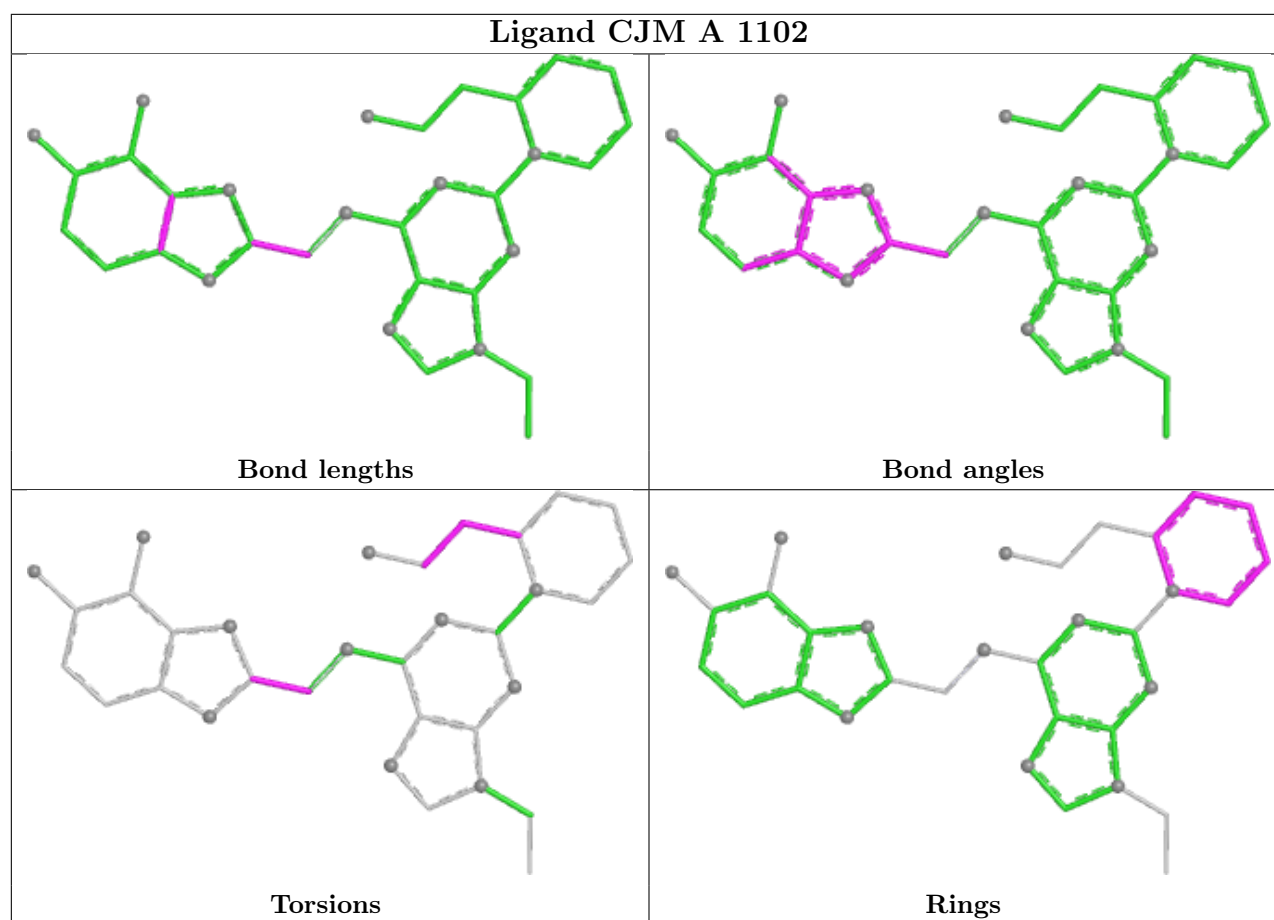
All (2) ring outliers are listed below:

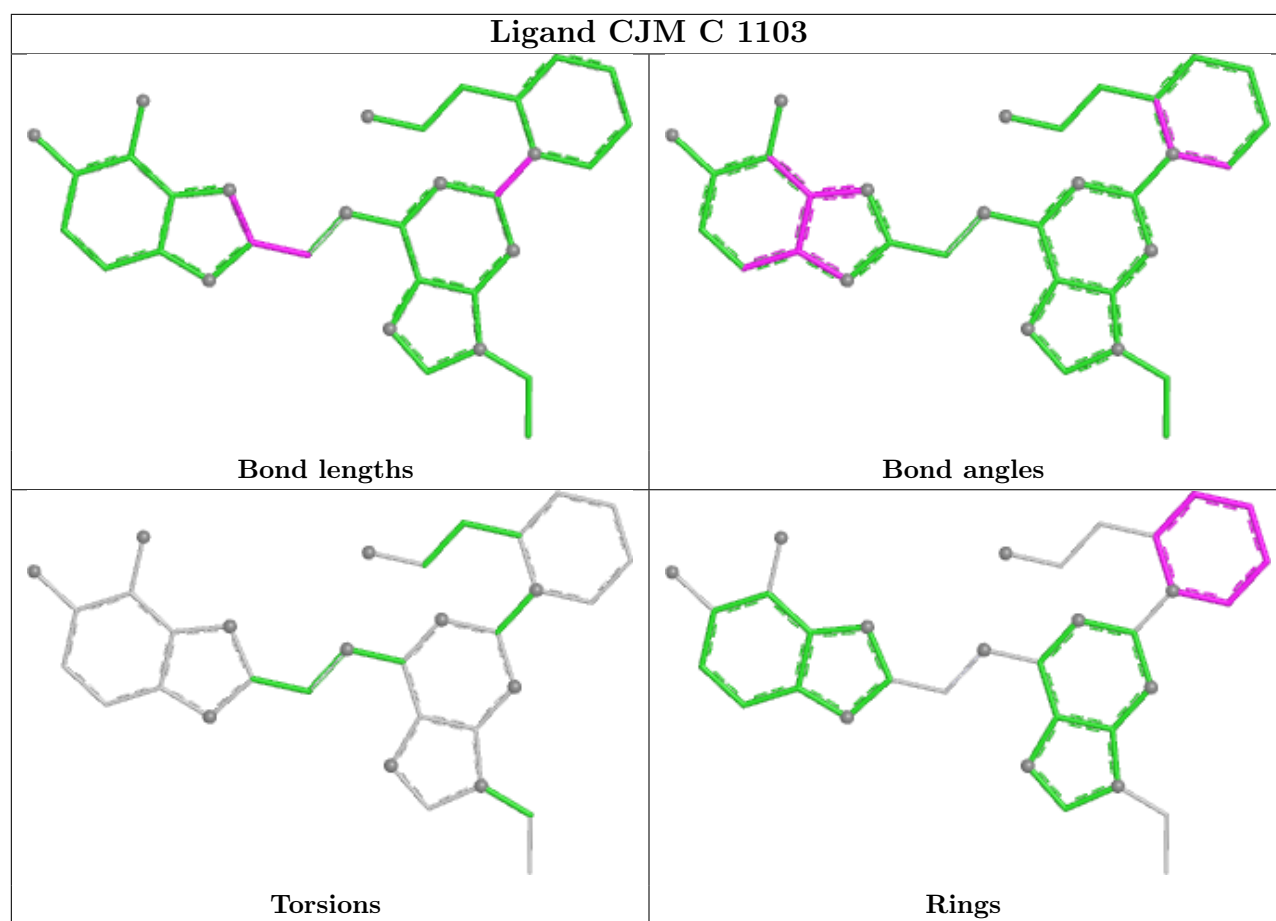
Mol	Chain	Res	Type	Atoms
4	A	1102	CJM	C26-C27-C28-C29-C30-N25
4	C	1103	CJM	C26-C27-C28-C29-C30-N25

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1102	CJM	4	0
5	B	301	EDO	1	0
5	C	1102	EDO	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 [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	309/320 (96%)	-0.21	1 (0%) 90 80	48, 83, 141, 169	0
1	C	307/320 (95%)	-0.08	1 (0%) 90 80	62, 99, 161, 198	0
2	B	240/267 (89%)	-0.51	0 100 100	41, 64, 113, 151	0
2	D	241/267 (90%)	-0.37	1 (0%) 88 75	44, 76, 138, 188	0
All	All	1097/1174 (93%)	-0.27	3 (0%) 90 80	41, 82, 147, 198	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1036	TRP	3.6
2	D	230	MET	2.6
1	C	738	TYR	2.4

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	TPO	A	893	11/12	0.88	0.11	93,96,101,102	0
1	TPO	C	893	11/12	0.95	0.07	83,85,89,91	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

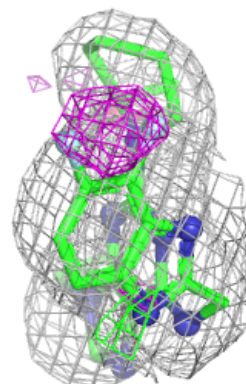
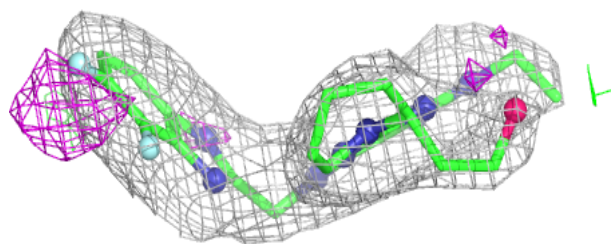
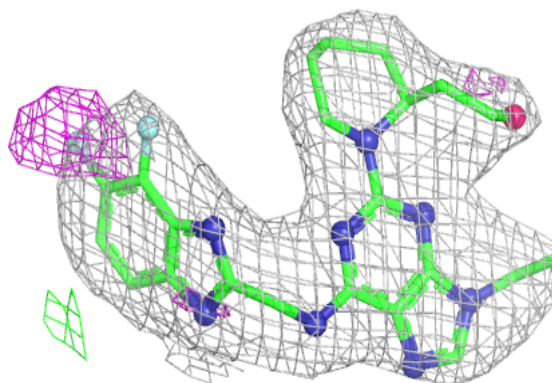
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	EDO	B	302	4/4	0.62	0.27	83,87,91,94	0
3	MG	C	1101	1/1	0.80	0.30	83,83,83,83	0
3	MG	A	1101	1/1	0.81	0.21	73,73,73,73	0
5	EDO	B	301	4/4	0.84	0.14	84,84,87,89	0
4	CJM	A	1102	33/33	0.93	0.09	53,59,69,71	0
4	CJM	C	1103	33/33	0.93	0.09	60,85,93,93	0
5	EDO	C	1102	4/4	0.94	0.10	48,52,58,62	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.

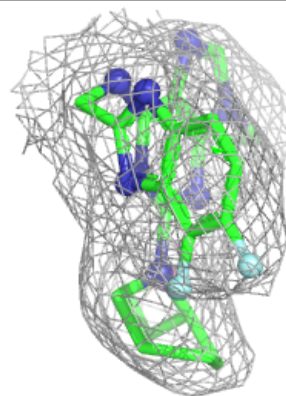
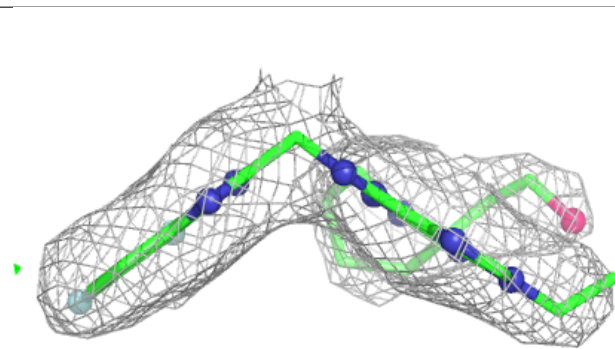
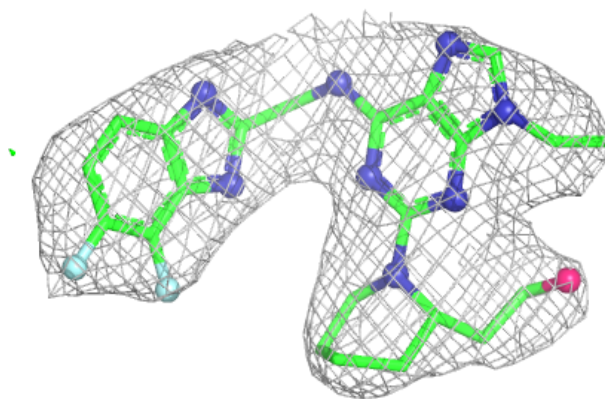
Electron density around CJM A 1102:

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 CJM C 1103:

$2mF_o-DF_c$ (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.