



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:39 PM UTC

PDB ID : 4R3U / pdb_00004r3u
Title : Crystal structure of 2-Hydroxyisobutyryl-CoA Mutase
Authors : Zahn, M.; Kurteva-Yaneva, N.; Rohwerder, T.; Straeter, N.
Deposited on : 2014-08-18
Resolution : 2.50 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

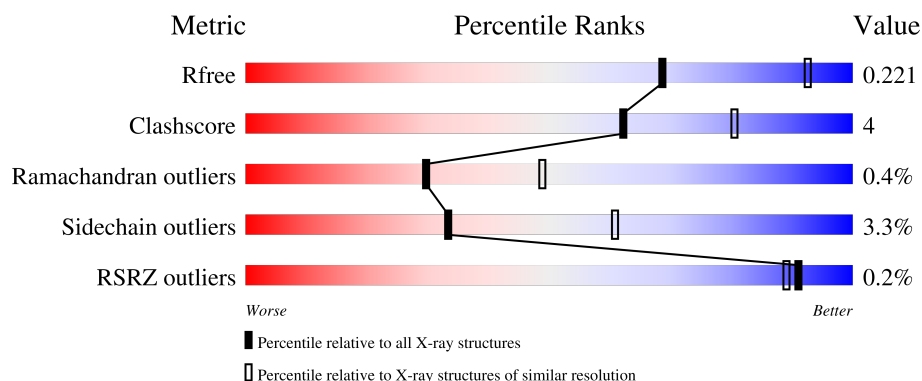
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	584	
1	B	584	
2	C	158	
2	D	158	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 11404 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 2-hydroxyisobutyryl-CoA mutase large subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	557	Total	C	N	O	S	0	0	0
			4423	2804	755	845	19			
1	B	555	Total	C	N	O	S	0	0	0
			4406	2795	753	839	19			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	MET	-	expression tag	UNP I3VE77
A	-9	ALA	-	expression tag	UNP I3VE77
A	-8	SER	-	expression tag	UNP I3VE77
A	-7	HIS	-	expression tag	UNP I3VE77
A	-6	HIS	-	expression tag	UNP I3VE77
A	-5	HIS	-	expression tag	UNP I3VE77
A	-4	HIS	-	expression tag	UNP I3VE77
A	-3	HIS	-	expression tag	UNP I3VE77
A	-2	HIS	-	expression tag	UNP I3VE77
A	-1	SER	-	expression tag	UNP I3VE77
A	0	GLY	-	expression tag	UNP I3VE77
A	563	GLY	-	expression tag	UNP I3VE77
A	564	SER	-	expression tag	UNP I3VE77
A	565	ALA	-	expression tag	UNP I3VE77
A	566	TRP	-	expression tag	UNP I3VE77
A	567	SER	-	expression tag	UNP I3VE77
A	568	HIS	-	expression tag	UNP I3VE77
A	569	PRO	-	expression tag	UNP I3VE77
A	570	GLN	-	expression tag	UNP I3VE77
A	571	PHE	-	expression tag	UNP I3VE77
A	572	GLU	-	expression tag	UNP I3VE77
A	573	LYS	-	expression tag	UNP I3VE77
B	-10	MET	-	expression tag	UNP I3VE77
B	-9	ALA	-	expression tag	UNP I3VE77
B	-8	SER	-	expression tag	UNP I3VE77

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	HIS	-	expression tag	UNP I3VE77
B	-6	HIS	-	expression tag	UNP I3VE77
B	-5	HIS	-	expression tag	UNP I3VE77
B	-4	HIS	-	expression tag	UNP I3VE77
B	-3	HIS	-	expression tag	UNP I3VE77
B	-2	HIS	-	expression tag	UNP I3VE77
B	-1	SER	-	expression tag	UNP I3VE77
B	0	GLY	-	expression tag	UNP I3VE77
B	563	GLY	-	expression tag	UNP I3VE77
B	564	SER	-	expression tag	UNP I3VE77
B	565	ALA	-	expression tag	UNP I3VE77
B	566	TRP	-	expression tag	UNP I3VE77
B	567	SER	-	expression tag	UNP I3VE77
B	568	HIS	-	expression tag	UNP I3VE77
B	569	PRO	-	expression tag	UNP I3VE77
B	570	GLN	-	expression tag	UNP I3VE77
B	571	PHE	-	expression tag	UNP I3VE77
B	572	GLU	-	expression tag	UNP I3VE77
B	573	LYS	-	expression tag	UNP I3VE77

- Molecule 2 is a protein called 2-hydroxyisobutyryl-CoA mutase small subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	131	Total	C	N	O	S	0	0	0
			982	620	176	184	2			
2	D	129	Total	C	N	O	S	0	0	0
			962	609	170	181	2			

There are 44 discrepancies between the modelled and reference sequences:

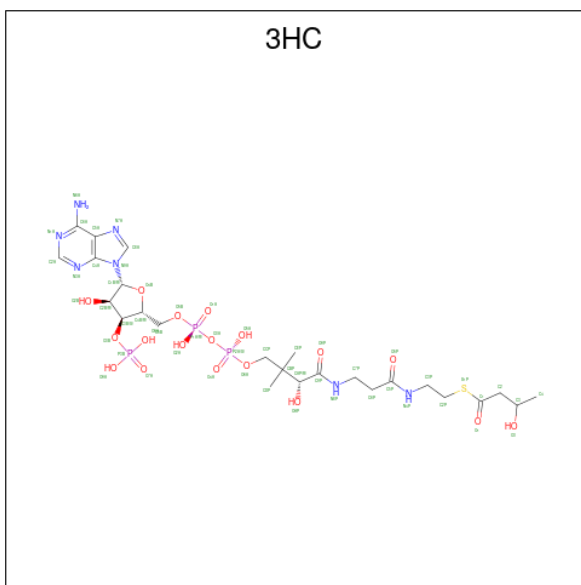
Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	MET	-	expression tag	UNP I3VE74
C	-9	ALA	-	expression tag	UNP I3VE74
C	-8	SER	-	expression tag	UNP I3VE74
C	-7	HIS	-	expression tag	UNP I3VE74
C	-6	HIS	-	expression tag	UNP I3VE74
C	-5	HIS	-	expression tag	UNP I3VE74
C	-4	HIS	-	expression tag	UNP I3VE74
C	-3	HIS	-	expression tag	UNP I3VE74
C	-2	HIS	-	expression tag	UNP I3VE74
C	-1	SER	-	expression tag	UNP I3VE74
C	0	GLY	-	expression tag	UNP I3VE74

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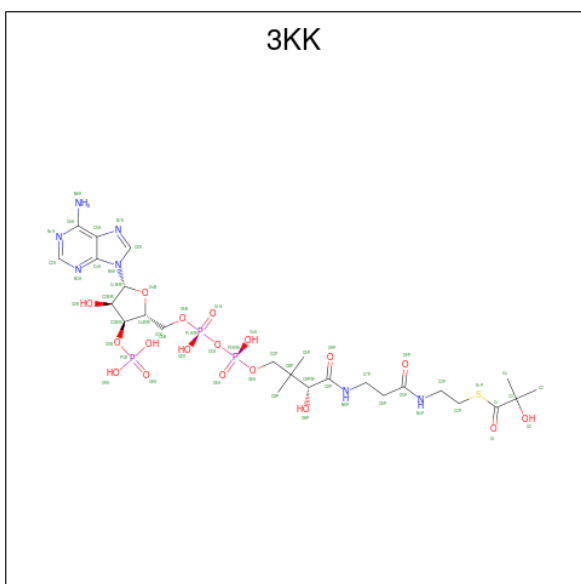
Chain	Residue	Modelled	Actual	Comment	Reference
C	137	GLY	-	expression tag	UNP I3VE74
C	138	SER	-	expression tag	UNP I3VE74
C	139	ALA	-	expression tag	UNP I3VE74
C	140	TRP	-	expression tag	UNP I3VE74
C	141	SER	-	expression tag	UNP I3VE74
C	142	HIS	-	expression tag	UNP I3VE74
C	143	PRO	-	expression tag	UNP I3VE74
C	144	GLN	-	expression tag	UNP I3VE74
C	145	PHE	-	expression tag	UNP I3VE74
C	146	GLU	-	expression tag	UNP I3VE74
C	147	LYS	-	expression tag	UNP I3VE74
D	-10	MET	-	expression tag	UNP I3VE74
D	-9	ALA	-	expression tag	UNP I3VE74
D	-8	SER	-	expression tag	UNP I3VE74
D	-7	HIS	-	expression tag	UNP I3VE74
D	-6	HIS	-	expression tag	UNP I3VE74
D	-5	HIS	-	expression tag	UNP I3VE74
D	-4	HIS	-	expression tag	UNP I3VE74
D	-3	HIS	-	expression tag	UNP I3VE74
D	-2	HIS	-	expression tag	UNP I3VE74
D	-1	SER	-	expression tag	UNP I3VE74
D	0	GLY	-	expression tag	UNP I3VE74
D	137	GLY	-	expression tag	UNP I3VE74
D	138	SER	-	expression tag	UNP I3VE74
D	139	ALA	-	expression tag	UNP I3VE74
D	140	TRP	-	expression tag	UNP I3VE74
D	141	SER	-	expression tag	UNP I3VE74
D	142	HIS	-	expression tag	UNP I3VE74
D	143	PRO	-	expression tag	UNP I3VE74
D	144	GLN	-	expression tag	UNP I3VE74
D	145	PHE	-	expression tag	UNP I3VE74
D	146	GLU	-	expression tag	UNP I3VE74
D	147	LYS	-	expression tag	UNP I3VE74

- Molecule 3 is 3-HYDROXYBUTANOYL-COENZYME A (CCD ID: 3HC) (formula: C₂₅H₄₂N₇O₁₈P₃S).



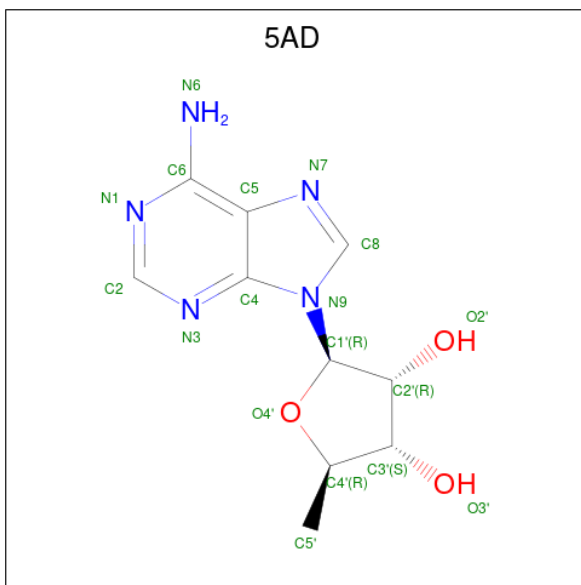
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	0	1
			54	25	7	18	3	1		
3	B	1	Total	C	N	O	P	S	0	1
			54	25	7	18	3	1		

- Molecule 4 is S-{(3R,5R,9R)-1-[(2R,3S,4R,5R)-5-(6-amino-9H-purin-9-yl)-4-hydroxy-3-(phosphonoxy)tetrahydrofuran-2-yl]-3,5,9-trihydroxy-8,8-dimethyl-3,5-dioxido-10,14-dioxo-2,4,6-trioxa-11,15-diaza-3lambda 5 ,5lambda 5 -diphosphaheptadecan-17-yl} 2-hydroxy-2-methylpropanethioate (CCD ID: 3KK) (formula: C₂₅H₄₂N₇O₁₈P₃S).



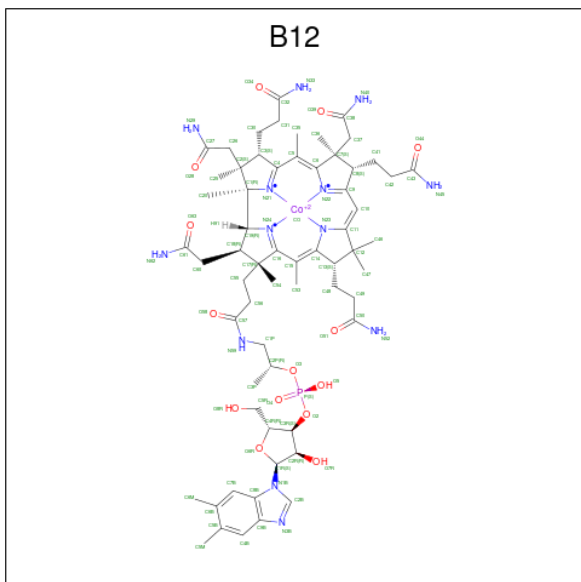
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	S	
			54	25	7	18	3	1	0
4	B	1	Total	C	N	O	P	S	
			54	25	7	18	3	1	0

- Molecule 5 is 5'-DEOXYADENOSINE (CCD ID: 5AD) (formula: $C_{10}H_{13}N_5O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O		
			18	10	5	3	0	1

- Molecule 6 is COBALAMIN (CCD ID: B12) (formula: $C_{62}H_{89}CoN_{13}O_{14}P$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
6	C	1	Total 91	C 62	Co 1	N 13	O 14	P 1	0	0
6	D	1	Total 91	C 62	Co 1	N 13	O 14	P 1	0	0

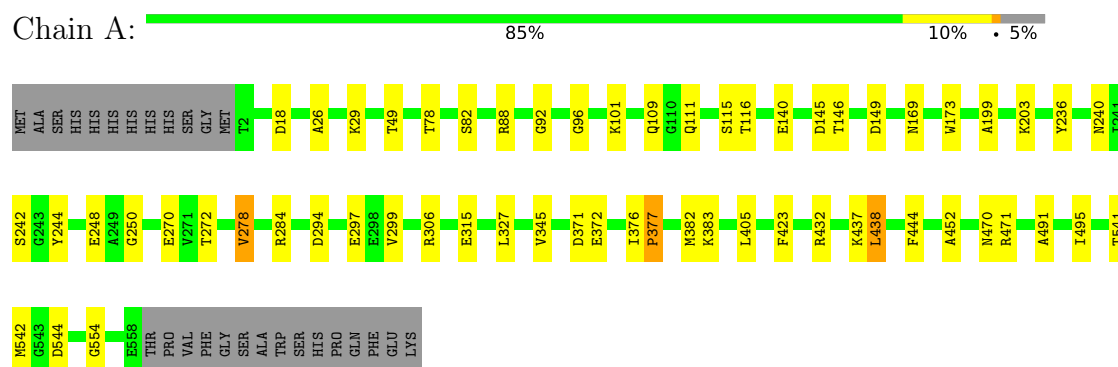
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	158	Total 158	O 158	0	5
7	B	16	Total 16	O 16	0	0
7	C	38	Total 38	O 38	0	1
7	D	3	Total 3	O 3	0	0

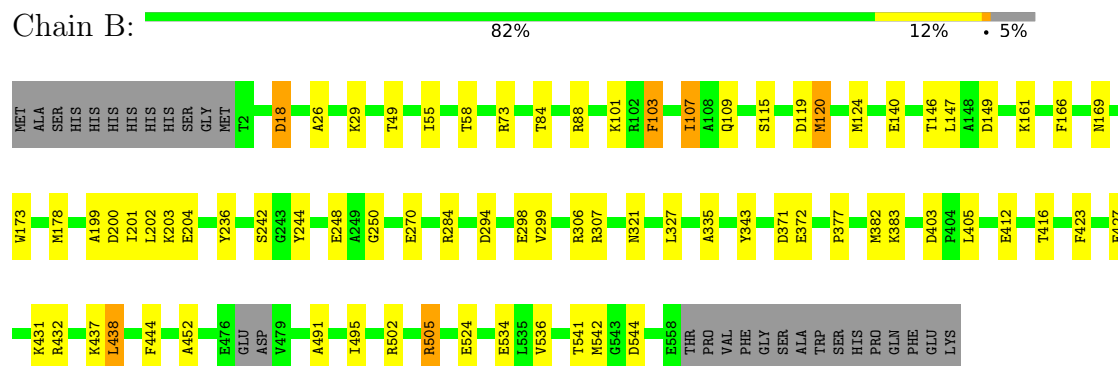
3 Residue-property plots

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

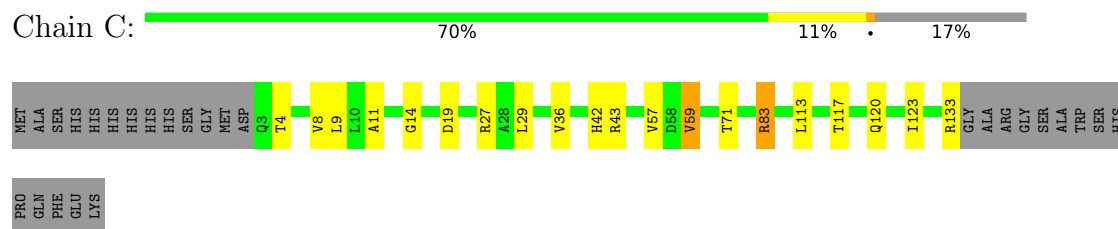
- Molecule 1: 2-hydroxyisobutyryl-CoA mutase large subunit



- Molecule 1: 2-hydroxyisobutyryl-CoA mutase large subunit



- Molecule 2: 2-hydroxyisobutyryl-CoA mutase small subunit



- Molecule 2: 2-hydroxyisobutyryl-CoA mutase small subunit



ALA
TRP
SER
HIS
PRO
GLN
PHE
GLU
LYS

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	69.03Å 119.56Å 173.94Å 90.00° 91.23° 90.00°	Depositor
Resolution (Å)	24.63 – 2.50 24.63 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (24.63-2.50) 99.9 (24.63-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.40 (at 2.50Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.168 , 0.217 0.172 , 0.221	Depositor DCC
R_{free} test set	1519 reflections (3.11%)	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.857	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 49.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.020 for -1/2*h+1/2*k,3/2*h+1/2*k,-l 0.017 for -1/2*h-1/2*k,-3/2*h+1/2*k,-l 0.010 for 1/2*h+1/2*k,3/2*h-1/2*k,-l 0.011 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.028 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11404	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.32% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 5AD, B12, 3HC, 3KK

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	0.89	0/4511	1.41	20/6100 (0.3%)
1	B	0.82	2/4493 (0.0%)	1.40	19/6074 (0.3%)
2	C	0.96	0/992	1.37	4/1346 (0.3%)
2	D	0.83	0/972	1.43	4/1320 (0.3%)
All	All	0.86	2/10968 (0.0%)	1.40	47/14840 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	120	MET	SD-CE	-7.79	1.60	1.79
1	B	124	MET	SD-CE	-5.35	1.66	1.79

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	371	ASP	CA-CB-CG	7.47	120.07	112.60
2	D	50	ASN	CA-CB-CG	7.25	119.85	112.60
1	B	377	PRO	N-CA-C	6.93	122.23	111.21
1	A	242	SER	N-CA-C	6.83	120.96	109.76
1	A	377	PRO	N-CA-C	6.75	122.11	111.38
1	B	119	ASP	CA-CB-CG	6.62	119.22	112.60
1	A	423	PHE	CA-C-N	6.50	129.00	120.28
1	A	423	PHE	C-N-CA	6.50	129.00	120.28
1	B	423	PHE	CA-C-N	6.50	128.99	120.28
1	B	423	PHE	C-N-CA	6.50	128.99	120.28
1	B	371	ASP	CA-CB-CG	6.31	118.91	112.60
1	B	103	PHE	CA-CB-CG	6.29	120.09	113.80
1	B	149	ASP	CA-CB-CG	6.29	118.89	112.60
1	A	149	ASP	CA-CB-CG	6.22	118.82	112.60
1	A	444	PHE	CA-C-N	6.13	128.49	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	444	PHE	C-N-CA	6.13	128.49	120.28
1	A	544	ASP	CA-C-N	5.82	127.89	120.56
1	A	544	ASP	C-N-CA	5.82	127.89	120.56
1	B	544	ASP	CA-C-N	5.79	127.86	120.56
1	B	544	ASP	C-N-CA	5.79	127.86	120.56
2	C	71	THR	CA-C-N	5.78	127.92	120.70
2	C	71	THR	C-N-CA	5.78	127.92	120.70
1	B	18	ASP	CA-C-N	5.73	127.95	120.28
1	B	18	ASP	C-N-CA	5.73	127.95	120.28
2	D	71	THR	CA-C-N	5.71	127.84	120.70
2	D	71	THR	C-N-CA	5.71	127.84	120.70
1	B	242	SER	N-CA-C	5.64	119.01	109.76
1	A	544	ASP	CA-CB-CG	5.64	118.24	112.60
1	B	166	PHE	CA-CB-CG	5.61	119.41	113.80
1	B	544	ASP	CA-CB-CG	5.54	118.14	112.60
2	C	19	ASP	CA-CB-CG	5.52	118.12	112.60
1	A	18	ASP	CA-C-N	5.51	127.66	120.28
1	A	18	ASP	C-N-CA	5.51	127.66	120.28
1	A	470	ASN	CA-CB-CG	5.46	118.06	112.60
2	D	11	ALA	N-CA-C	5.30	118.38	109.95
1	A	376	ILE	CB-CA-C	5.22	115.99	110.88
2	C	11	ALA	N-CA-C	5.21	118.09	109.85
1	B	542	MET	CA-C-N	5.11	125.65	119.98
1	B	542	MET	C-N-CA	5.11	125.65	119.98
1	B	444	PHE	CA-C-N	5.08	127.09	120.28
1	B	444	PHE	C-N-CA	5.08	127.09	120.28
1	A	96	GLY	CA-C-N	5.03	127.02	120.28
1	A	96	GLY	C-N-CA	5.03	127.02	120.28
1	B	321	ASN	CA-CB-CG	5.02	117.62	112.60
1	A	542	MET	CA-C-N	5.02	125.55	119.98
1	A	542	MET	C-N-CA	5.02	125.55	119.98
1	A	315	GLU	CB-CG-CD	5.01	121.12	112.60

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	4423	0	4391	29	0
1	B	4406	0	4380	34	0
2	C	982	0	1034	8	0
2	D	962	0	1013	9	0
3	A	54	0	38	4	0
3	B	54	0	38	0	0
4	A	54	0	38	3	0
4	B	54	0	38	0	0
5	A	18	0	13	6	0
6	C	91	0	88	9	0
6	D	91	0	88	9	0
7	A	158	0	0	3	0
7	B	16	0	0	1	0
7	C	38	0	0	0	0
7	D	3	0	0	0	0
All	All	11404	0	11159	92	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 4.

All (92) 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:B:307:ARG:HH22	1:B:416:THR:HG22	1.32	0.95
1:B:178:MET:HG2	1:B:536:VAL:HG21	1.49	0.94
1:B:178:MET:CG	1:B:536:VAL:HG21	2.01	0.90
6:C:800:B12:H531	6:C:800:B12:H552	1.60	0.83
1:A:78:THR:HG21	7:A:939:HOH:O	1.82	0.79
5:A:803[C]:5AD:H8	6:C:800:B12:H372	1.65	0.79
1:B:120:MET:HE1	6:D:800:B12:O44	1.85	0.77
6:D:800:B12:H552	6:D:800:B12:H531	1.64	0.77
1:A:491:ALA:O	1:A:495:ILE:HG12	1.90	0.70
1:B:307:ARG:NH2	1:B:416:THR:HG22	2.05	0.69
6:D:800:B12:H601	6:D:800:B12:H252	1.76	0.67
1:B:502:ARG:HG2	1:B:505:ARG:HH21	1.59	0.66
6:C:800:B12:H531	6:C:800:B12:C55	2.25	0.66
1:B:58:THR:OG1	1:B:416:THR:HG21	1.98	0.64
1:B:178:MET:HG2	1:B:536:VAL:CG2	2.25	0.64
6:D:800:B12:H531	6:D:800:B12:C55	2.29	0.63
1:B:178:MET:HG3	1:B:536:VAL:HG21	1.78	0.62
6:D:800:B12:H351	6:D:800:B12:H362	1.81	0.62
1:A:78:THR:HG23	4:A:802[B]:3KK:H42	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:77:PHE:CE1	2:D:90:VAL:HG11	2.36	0.61
6:C:800:B12:H601	6:C:800:B12:H252	1.81	0.61
6:C:800:B12:H362	6:C:800:B12:H351	1.82	0.60
1:A:92:GLY:O	5:A:803[C]:5AD:N6	2.34	0.60
2:D:77:PHE:CZ	2:D:90:VAL:HG11	2.36	0.60
5:A:803[C]:5AD:H5'3	7:A:978:HOH:O	2.01	0.59
1:A:78:THR:HG23	3:A:801[A]:3HC:H61A	1.66	0.59
1:B:299:VAL:HG11	1:B:405:LEU:HD12	1.84	0.59
1:B:147:LEU:HB2	1:B:536:VAL:HG22	1.84	0.58
1:A:306:ARG:HG2	1:A:327:LEU:HB3	1.85	0.58
1:A:299:VAL:HG11	1:A:405:LEU:HD12	1.85	0.57
1:B:306:ARG:HG2	1:B:327:LEU:HB3	1.86	0.57
1:B:412:GLU:O	1:B:416:THR:HG23	2.04	0.57
1:B:432:ARG:HB3	1:B:438:LEU:HD11	1.87	0.55
1:A:432:ARG:HB3	1:A:438:LEU:HD11	1.88	0.55
1:A:250:GLY:HA3	1:A:452:ALA:HB2	1.89	0.54
1:A:140:GLU:HB3	6:C:800:B12:H362	1.91	0.53
1:B:120:MET:HE1	6:D:800:B12:C43	2.39	0.52
1:B:73:ARG:HD2	1:B:403:ASP:OD2	2.08	0.52
1:B:140:GLU:HB3	6:D:800:B12:H362	1.93	0.51
1:B:502:ARG:HG2	1:B:505:ARG:NH2	2.24	0.51
1:B:250:GLY:HA3	1:B:452:ALA:HB2	1.93	0.51
2:C:83:ARG:HG2	2:C:83:ARG:HH21	1.76	0.50
1:B:201:ILE:O	1:B:204:GLU:HB2	2.12	0.50
1:B:491:ALA:O	1:B:495:ILE:HG12	2.13	0.49
1:A:145:ASP:OD1	7:A:1003:HOH:O	2.20	0.48
1:A:372:GLU:OE1	5:A:803[C]:5AD:H3'	2.14	0.48
1:B:382:MET:HE2	1:B:382:MET:HA	1.95	0.48
5:A:803[C]:5AD:C8	6:C:800:B12:H372	2.41	0.47
1:B:55:ILE:HA	7:B:913:HOH:O	2.14	0.47
1:B:88:ARG:HG2	1:B:115:SER:HB2	1.95	0.47
2:D:14:GLY:O	2:D:42:HIS:HA	2.14	0.47
1:A:146:THR:HG22	1:A:541:THR:HG22	1.96	0.47
1:A:377:PRO:HD2	5:A:803[C]:5AD:C2	2.45	0.47
2:D:29:LEU:HD23	2:D:123:ILE:HD13	1.96	0.47
2:C:14:GLY:O	2:C:42:HIS:HA	2.15	0.46
1:A:88:ARG:HG2	1:A:115:SER:HB2	1.97	0.46
1:A:294:ASP:HB3	1:A:297:GLU:HB3	1.98	0.46
1:A:240:ASN:HD21	4:A:802[B]:3KK:H14	1.81	0.45
1:B:103:PHE:O	1:B:107:ILE:HG23	2.16	0.45
2:C:8:VAL:HG13	2:C:59:VAL:HG22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:GLN:HG3	1:A:383:LYS:HG3	1.97	0.45
1:A:382:MET:HA	1:A:382:MET:HE2	1.98	0.45
1:A:240:ASN:HD21	3:A:801[A]:3HC:H61	1.82	0.44
2:C:83:ARG:HH21	2:C:83:ARG:CG	2.30	0.44
1:A:26:ALA:HA	1:A:29:LYS:HE3	1.98	0.44
1:B:109:GLN:HG3	1:B:383:LYS:HG3	2.00	0.44
2:D:119:PRO:O	2:D:123:ILE:HG12	2.17	0.44
1:A:272:THR:HG21	1:A:278:VAL:HG22	2.00	0.44
2:D:59:VAL:HG22	2:D:89:ILE:HB	2.01	0.43
1:A:236:TYR:O	1:A:284:ARG:HD2	2.17	0.43
1:B:26:ALA:HA	1:B:29:LYS:HE3	2.00	0.43
2:C:120:GLN:HA	2:C:123:ILE:HD12	1.99	0.43
1:A:554:GLY:HA2	1:B:431:LYS:O	2.18	0.43
1:B:236:TYR:O	1:B:284:ARG:HD2	2.18	0.42
2:D:119:PRO:HA	2:D:122:ILE:HD12	2.01	0.42
1:B:146:THR:HG22	1:B:541:THR:HG22	2.00	0.42
1:B:244:TYR:CE2	1:B:248:GLU:HG3	2.55	0.42
1:A:78:THR:OG1	1:A:82:SER:HB2	2.20	0.42
1:A:240:ASN:ND2	4:A:802[B]:3KK:H14	2.35	0.41
2:C:9:LEU:HB2	2:C:57:VAL:HG11	2.02	0.41
1:A:248:GLU:HB3	6:C:800:B12:H532	2.02	0.41
2:D:4:THR:HA	2:D:5:PRO:HD3	1.92	0.41
1:A:244:TYR:CE2	1:A:248:GLU:HG3	2.55	0.41
3:A:801[A]:3HC:HC41	6:C:800:B12:H371	2.02	0.41
1:B:294:ASP:O	1:B:298:GLU:HG2	2.21	0.41
2:C:113:LEU:HB3	2:C:117:THR:HG21	2.01	0.41
2:D:29:LEU:HB2	2:D:36:VAL:HG21	2.03	0.41
1:B:248:GLU:HB3	6:D:800:B12:H532	2.02	0.40
1:A:240:ASN:ND2	3:A:801[A]:3HC:H61	2.36	0.40
2:C:29:LEU:HB2	2:C:36:VAL:HG21	2.02	0.40
1:B:335:ALA:HB1	1:B:372:GLU:OE2	2.22	0.40
6:D:800:B12:H362	6:D:800:B12:C35	2.51	0.40

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	555/584 (95%)	533 (96%)	20 (4%)	2 (0%)	30	49
1	B	551/584 (94%)	531 (96%)	16 (3%)	4 (1%)	18	34
2	C	129/158 (82%)	125 (97%)	4 (3%)	0	100	100
2	D	127/158 (80%)	123 (97%)	4 (3%)	0	100	100
All	All	1362/1484 (92%)	1312 (96%)	44 (3%)	6 (0%)	30	49

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	169	ASN
1	B	169	ASN
1	A	199	ALA
1	B	199	ALA
1	B	343	TYR
1	B	200	ASP

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	467/490 (95%)	455 (97%)	12 (3%)	40	68
1	B	465/490 (95%)	449 (97%)	16 (3%)	32	60
2	C	106/127 (84%)	100 (94%)	6 (6%)	18	39
2	D	104/127 (82%)	100 (96%)	4 (4%)	29	56
All	All	1142/1234 (92%)	1104 (97%)	38 (3%)	33	61

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	THR
1	A	101	LYS
1	A	111	GLN
1	A	116	THR
1	A	173	TRP
1	A	203	LYS
1	A	270	GLU
1	A	278	VAL
1	A	345	VAL
1	A	437	LYS
1	A	438	LEU
1	A	471	ARG
1	B	18	ASP
1	B	49	THR
1	B	84	THR
1	B	101	LYS
1	B	107	ILE
1	B	161	LYS
1	B	173	TRP
1	B	202	LEU
1	B	203	LYS
1	B	270	GLU
1	B	427	GLU
1	B	437	LYS
1	B	438	LEU
1	B	505	ARG
1	B	524	GLU
1	B	534	GLU
2	C	4	THR
2	C	27	ARG
2	C	43	ARG
2	C	59	VAL
2	C	83	ARG
2	C	133	ARG
2	D	62	VAL
2	D	104	ILE
2	D	106	LYS
2	D	111	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN

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Mol	Chain	Res	Type
1	A	47	GLN
1	A	240	ASN
1	A	342	GLN
1	A	441	GLN
1	A	526	GLN
1	B	13	GLN
1	B	47	GLN
1	B	198	GLN
1	B	240	ASN
1	B	441	GLN
1	B	526	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 ⓘ

7 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	$\# Z > 2$
6	B12	C	800	2	94,101,101	0.92	3 (3%)	149,166,166	1.08	9 (6%)
5	5AD	A	803[C]	-	20,20,20	1.78	6 (30%)	28,30,30	2.57	12 (42%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	3KK	B	802[B]	-	53,56,56	1.59	9 (16%)	75,85,85	1.67	12 (16%)
6	B12	D	800	2	94,101,101	0.96	4 (4%)	149,166,166	1.06	9 (6%)
3	3HC	B	801[A]	-	54,56,56	1.97	7 (12%)	76,83,83	1.88	13 (17%)
4	3KK	A	802[B]	-	53,56,56	1.64	8 (15%)	75,85,85	1.67	12 (16%)
3	3HC	A	801[A]	-	54,56,56	2.01	7 (12%)	76,83,83	1.90	13 (17%)

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	B12	C	800	2	-	11/56/223/223	0/3/11/11
5	5AD	A	803[C]	-	-	0/4/20/20	0/3/3/3
4	3KK	B	802[B]	-	-	2/56/73/73	0/3/3/3
6	B12	D	800	2	-	12/56/223/223	0/3/11/11
3	3HC	B	801[A]	-	-	5/55/71/71	0/3/3/3
4	3KK	A	802[B]	-	-	3/56/73/73	0/3/3/3
3	3HC	A	801[A]	-	-	6/55/71/71	0/3/3/3

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801[A]	3HC	C2A-N1A	8.24	1.48	1.33
3	B	801[A]	3HC	C2A-N1A	8.11	1.48	1.33
3	B	801[A]	3HC	C2A-N3A	7.60	1.47	1.33
3	A	801[A]	3HC	C2A-N3A	7.54	1.47	1.33
6	D	800	B12	C14-N23	6.46	1.43	1.35
6	C	800	B12	C14-N23	5.77	1.42	1.35
4	A	802[B]	3KK	C2A-N1A	4.94	1.42	1.33
4	B	802[B]	3KK	C2A-N1A	4.90	1.42	1.33
4	A	802[B]	3KK	P1A-O3A	4.83	1.64	1.59
3	A	801[A]	3HC	P1A-O3A	4.75	1.64	1.59
5	A	803[C]	5AD	C2-N3	4.67	1.42	1.33
4	A	802[B]	3KK	C2A-N3A	4.41	1.41	1.33
4	B	802[B]	3KK	C2A-N3A	4.40	1.41	1.33
4	B	802[B]	3KK	P1A-O3A	4.22	1.64	1.59
3	B	801[A]	3HC	P1A-O3A	4.18	1.64	1.59
3	B	801[A]	3HC	P2A-O3A	3.55	1.63	1.59
4	B	802[B]	3KK	P2A-O3A	3.52	1.63	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	803[C]	5AD	C2-N1	3.35	1.39	1.33
4	B	802[B]	3KK	O1-C1	3.29	1.23	1.20
4	A	802[B]	3KK	O1-C1	3.27	1.23	1.20
3	A	801[A]	3HC	P2A-O3A	3.20	1.62	1.59
6	C	800	B12	C9-N22	3.19	1.38	1.30
4	A	802[B]	3KK	P2A-O3A	3.08	1.62	1.59
6	C	800	B12	C1P-C2P	2.89	1.58	1.51
3	A	801[A]	3HC	C5A-N7A	-2.76	1.34	1.39
6	D	800	B12	C9-N22	2.70	1.37	1.30
3	B	801[A]	3HC	C5A-N7A	-2.68	1.34	1.39
4	A	802[B]	3KK	C5A-N7A	-2.66	1.34	1.39
5	A	803[C]	5AD	C5-N7	-2.66	1.34	1.39
6	D	800	B12	C1P-C2P	2.62	1.58	1.51
4	B	802[B]	3KK	C5A-N7A	-2.60	1.34	1.39
3	A	801[A]	3HC	C8A-N9A	-2.44	1.33	1.37
4	A	802[B]	3KK	C8A-N9A	-2.31	1.33	1.37
3	B	801[A]	3HC	C8A-N9A	-2.29	1.33	1.37
5	A	803[C]	5AD	C5-C6	-2.28	1.34	1.41
6	D	800	B12	C16-C15	2.24	1.51	1.44
4	B	802[B]	3KK	C8A-N9A	-2.20	1.33	1.37
3	B	801[A]	3HC	C5A-C4A	-2.19	1.35	1.39
4	B	802[B]	3KK	C5A-C4A	-2.18	1.35	1.39
4	A	802[B]	3KK	C5A-C4A	-2.11	1.35	1.39
5	A	803[C]	5AD	C8-N9	-2.05	1.34	1.37
3	A	801[A]	3HC	C5A-C4A	-2.04	1.35	1.39
5	A	803[C]	5AD	C5-C4	-2.04	1.35	1.39
4	B	802[B]	3KK	C5A-C6A	-2.00	1.35	1.41

All (80) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	801[A]	3HC	N3A-C2A-N1A	-10.81	112.22	128.58
3	A	801[A]	3HC	N3A-C2A-N1A	-10.81	112.22	128.58
4	B	802[B]	3KK	N1A-C2A-N3A	-8.15	116.25	128.58
4	A	802[B]	3KK	N1A-C2A-N3A	-8.13	116.27	128.58
5	A	803[C]	5AD	N3-C2-N1	-6.37	118.94	128.58
5	A	803[C]	5AD	C5'-C4'-C3'	-5.64	109.77	115.70
5	A	803[C]	5AD	C5-C4-N3	-4.71	120.24	126.72
3	B	801[A]	3HC	C2A-N3A-C4A	4.08	121.80	111.83
4	B	802[B]	3KK	N9A-C8A-N7A	-4.04	108.20	113.94
4	A	802[B]	3KK	N9A-C8A-N7A	-4.03	108.21	113.94
3	A	801[A]	3HC	C2A-N3A-C4A	4.01	121.63	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	801[A]	3HC	N9A-C8A-N7A	-3.97	108.31	113.94
3	A	801[A]	3HC	N9A-C8A-N7A	-3.96	108.32	113.94
4	B	802[B]	3KK	C5A-C4A-N3A	-3.92	121.31	126.72
4	A	802[B]	3KK	C5A-C4A-N3A	-3.89	121.36	126.72
4	B	802[B]	3KK	C2A-N3A-C4A	3.57	120.56	111.83
6	C	800	B12	C2-C1-C19	-3.55	113.09	118.61
4	A	802[B]	3KK	C2A-N3A-C4A	3.49	120.35	111.83
6	C	800	B12	C9-N22-C6	3.46	109.44	105.28
3	B	801[A]	3HC	C5A-C4A-N3A	-3.45	121.96	126.72
3	A	801[A]	3HC	C5A-C4A-N3A	-3.41	122.02	126.72
6	D	800	B12	C2-C1-C19	-3.30	113.48	118.61
6	D	800	B12	C20-C1-C2	3.27	118.67	113.28
6	D	800	B12	C9-N22-C6	3.21	109.14	105.28
6	C	800	B12	C20-C1-C2	3.19	118.55	113.28
5	A	803[C]	5AD	C1'-N9-C8	-3.19	120.02	127.09
5	A	803[C]	5AD	C5-C6-N6	-3.17	115.43	123.29
5	A	803[C]	5AD	N9-C8-N7	-3.17	109.44	113.94
4	A	802[B]	3KK	C5A-N7A-C8A	3.09	108.31	103.45
5	A	803[C]	5AD	N3-C4-N9	3.09	132.42	127.17
3	A	801[A]	3HC	C5A-N7A-C8A	3.09	108.31	103.45
4	B	802[B]	3KK	C5A-N7A-C8A	3.06	108.26	103.45
5	A	803[C]	5AD	C2-N3-C4	3.04	119.25	111.83
3	B	801[A]	3HC	C5A-N7A-C8A	3.02	108.20	103.45
3	A	801[A]	3HC	C4A-C5A-N7A	-2.91	107.26	110.58
4	A	802[B]	3KK	C4A-C5A-N7A	-2.80	107.38	110.58
3	B	801[A]	3HC	C4A-C5A-N7A	-2.75	107.44	110.58
5	A	803[C]	5AD	C2'-C3'-C4'	2.71	106.36	102.36
6	C	800	B12	C10-C9-N22	2.68	128.79	125.74
4	B	802[B]	3KK	C4A-C5A-N7A	-2.67	107.52	110.58
4	B	802[B]	3KK	O1-C1-S1P	2.67	124.64	121.98
3	A	801[A]	3HC	O1-C1-C2	-2.60	119.66	123.74
3	B	801[A]	3HC	O1-C1-C2	-2.58	119.68	123.74
6	D	800	B12	C19-N24-C16	2.50	110.02	107.29
6	C	800	B12	C19-N24-C16	2.49	110.02	107.29
5	A	803[C]	5AD	C6-C5-C4	2.45	120.52	117.18
4	A	802[B]	3KK	O1-C1-S1P	2.44	124.41	121.98
4	B	802[B]	3KK	C4A-N9A-C8A	2.40	108.25	105.74
4	B	802[B]	3KK	C5A-C6A-N6A	-2.40	117.35	123.29
4	A	802[B]	3KK	C4A-N9A-C8A	2.39	108.25	105.74
5	A	803[C]	5AD	C5-N7-C8	2.37	107.17	103.45
6	D	800	B12	C10-C9-N22	2.36	128.44	125.74
3	B	801[A]	3HC	C5A-C6A-N6A	-2.35	117.46	123.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801[A]	3HC	C6A-C5A-C4A	2.33	120.36	117.18
3	A	801[A]	3HC	C2-C1-S1P	2.32	116.49	113.56
4	A	802[B]	3KK	C5A-C6A-N6A	-2.31	117.57	123.29
3	A	801[A]	3HC	C2P-C3P-N4P	-2.30	107.61	112.41
3	A	801[A]	3HC	C4A-N9A-C8A	2.30	108.15	105.74
3	B	801[A]	3HC	C2P-C3P-N4P	-2.30	107.61	112.41
3	A	801[A]	3HC	C5A-C6A-N6A	-2.27	117.65	123.29
6	C	800	B12	P-O3-C2P	2.27	128.86	121.10
3	B	801[A]	3HC	C4A-N9A-C8A	2.27	108.12	105.74
6	D	800	B12	C60-C18-C17	2.24	121.95	115.82
6	D	800	B12	O3-P-O4	2.22	116.91	109.81
4	B	802[B]	3KK	N3A-C4A-N9A	2.22	130.94	127.17
6	C	800	B12	C13-C12-C11	-2.21	98.50	100.97
3	B	801[A]	3HC	C6A-C5A-C4A	2.20	120.19	117.18
6	C	800	B12	O3-P-O4	2.18	116.78	109.81
6	D	800	B12	C54-C17-C55	-2.18	105.64	109.27
4	A	802[B]	3KK	C6A-C5A-C4A	2.18	120.15	117.18
3	B	801[A]	3HC	C2-C1-S1P	2.15	116.27	113.56
4	A	802[B]	3KK	N3A-C4A-N9A	2.14	130.81	127.17
4	A	802[B]	3KK	C2P-C3P-N4P	-2.14	107.94	112.41
4	B	802[B]	3KK	C2P-C3P-N4P	-2.14	107.94	112.41
6	D	800	B12	C8-C9-C10	-2.11	118.83	123.33
6	C	800	B12	C35-C5-C6	2.08	125.77	122.41
3	B	801[A]	3HC	O9A-P3B-O8A	2.07	115.56	107.80
5	A	803[C]	5AD	C5-C6-N1	2.07	122.76	117.51
3	A	801[A]	3HC	O9A-P3B-O8A	2.05	115.50	107.80
4	B	802[B]	3KK	C6A-C5A-C4A	2.04	119.96	117.18

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	801[A]	3HC	O1-C1-S1P-C2P
3	A	801[A]	3HC	C2-C1-S1P-C2P
3	B	801[A]	3HC	O1-C1-S1P-C2P
3	B	801[A]	3HC	C2-C1-S1P-C2P
6	C	800	B12	C2R-C1R-N1B-C8B
6	D	800	B12	C2R-C1R-N1B-C8B
6	D	800	B12	O6R-C1R-N1B-C8B
6	C	800	B12	O6R-C1R-N1B-C8B
6	C	800	B12	C2R-C1R-N1B-C2B
4	A	802[B]	3KK	O9P-C9P-CAP-OAP

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Mol	Chain	Res	Type	Atoms
4	B	802[B]	3KK	O9P-C9P-CAP-OAP
6	C	800	B12	C3-C30-C31-C32
6	D	800	B12	C3-C30-C31-C32
6	D	800	B12	C2R-C1R-N1B-C2B
6	D	800	B12	C55-C56-C57-O58
3	A	801[A]	3HC	O9P-C9P-CAP-OAP
3	B	801[A]	3HC	O9P-C9P-CAP-OAP
6	C	800	B12	C55-C56-C57-O58
6	D	800	B12	C55-C56-C57-N59
6	C	800	B12	C42-C41-C8-C9
6	D	800	B12	O6R-C1R-N1B-C2B
6	C	800	B12	C55-C56-C57-N59
6	C	800	B12	O6R-C1R-N1B-C2B
6	D	800	B12	C42-C41-C8-C9
6	C	800	B12	C2P-O3-P-O4
6	C	800	B12	C2P-O3-P-O5
6	D	800	B12	C2P-O3-P-O4
6	D	800	B12	C2P-O3-P-O5
3	A	801[A]	3HC	S1P-C1-C2-C3
3	B	801[A]	3HC	S1P-C1-C2-C3
3	A	801[A]	3HC	C4B-C5B-O5B-P1A
6	C	800	B12	C30-C31-C32-N33
6	D	800	B12	C30-C31-C32-O34
6	D	800	B12	C30-C31-C32-N33
3	A	801[A]	3HC	N8P-C9P-CAP-OAP
3	B	801[A]	3HC	N8P-C9P-CAP-OAP
4	A	802[B]	3KK	N8P-C9P-CAP-OAP
4	B	802[B]	3KK	N8P-C9P-CAP-OAP
4	A	802[B]	3KK	C4B-C5B-O5B-P1A

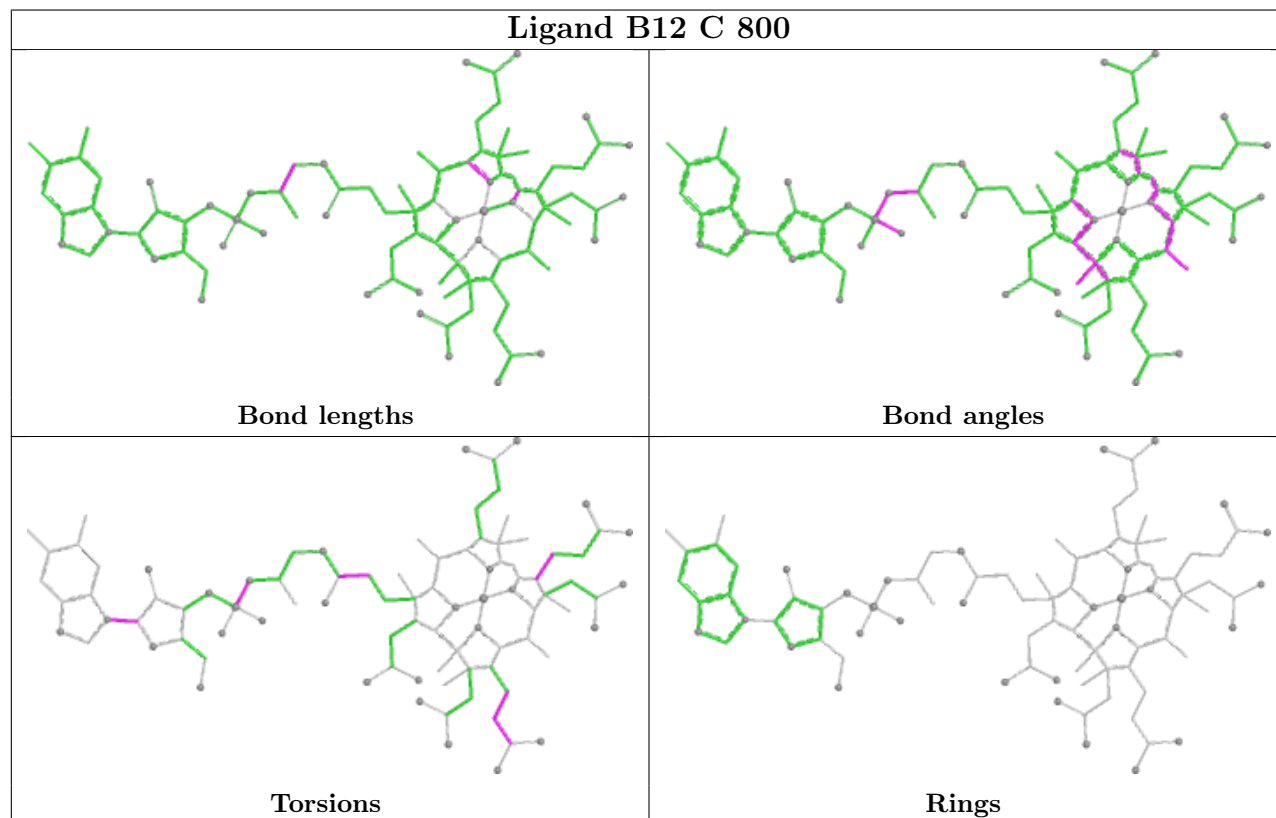
There are no ring outliers.

5 monomers are involved in 28 short contacts:

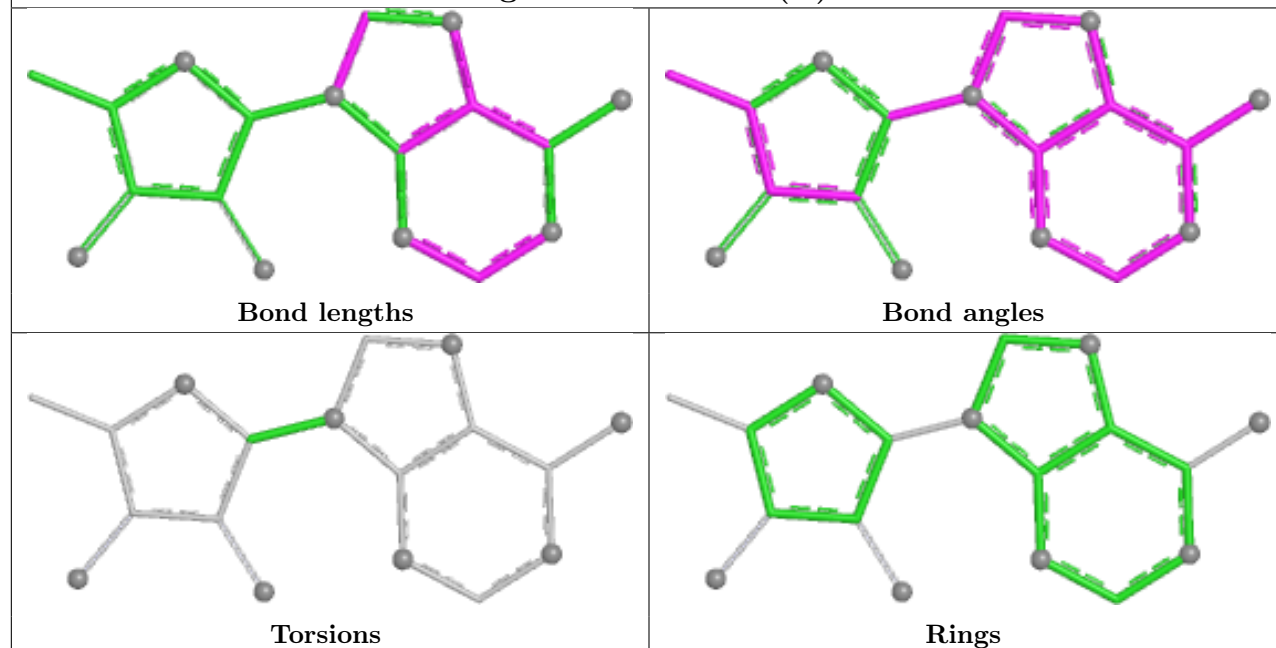
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	800	B12	9	0
5	A	803[C]	5AD	6	0
6	D	800	B12	9	0
4	A	802[B]	3KK	3	0
3	A	801[A]	3HC	4	0

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

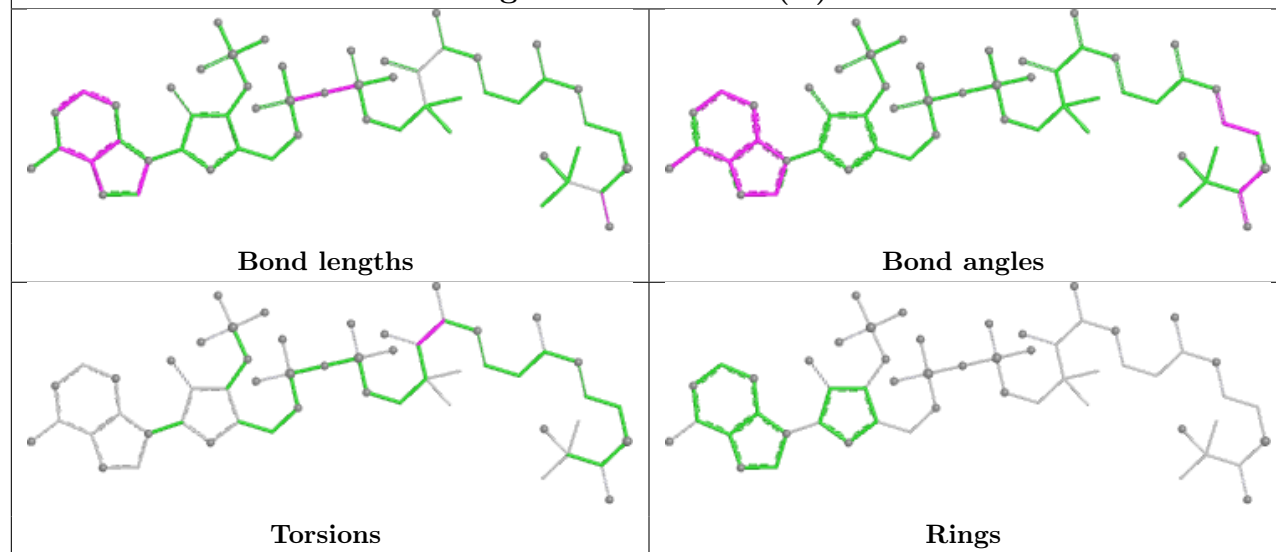
addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. 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.

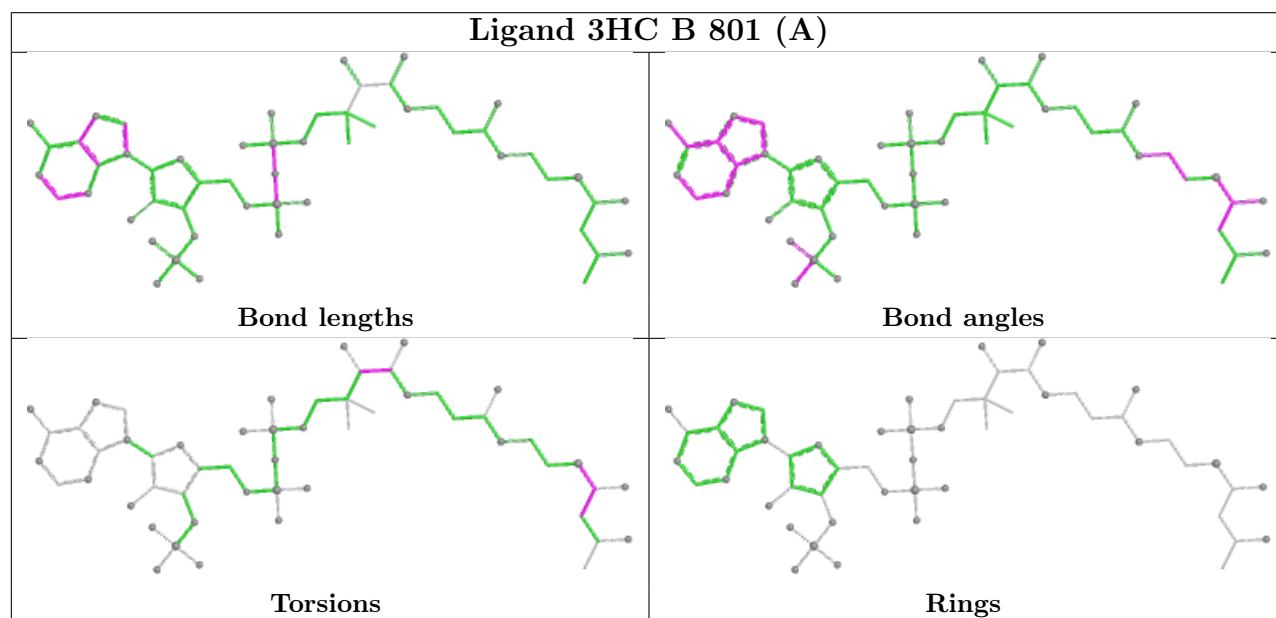
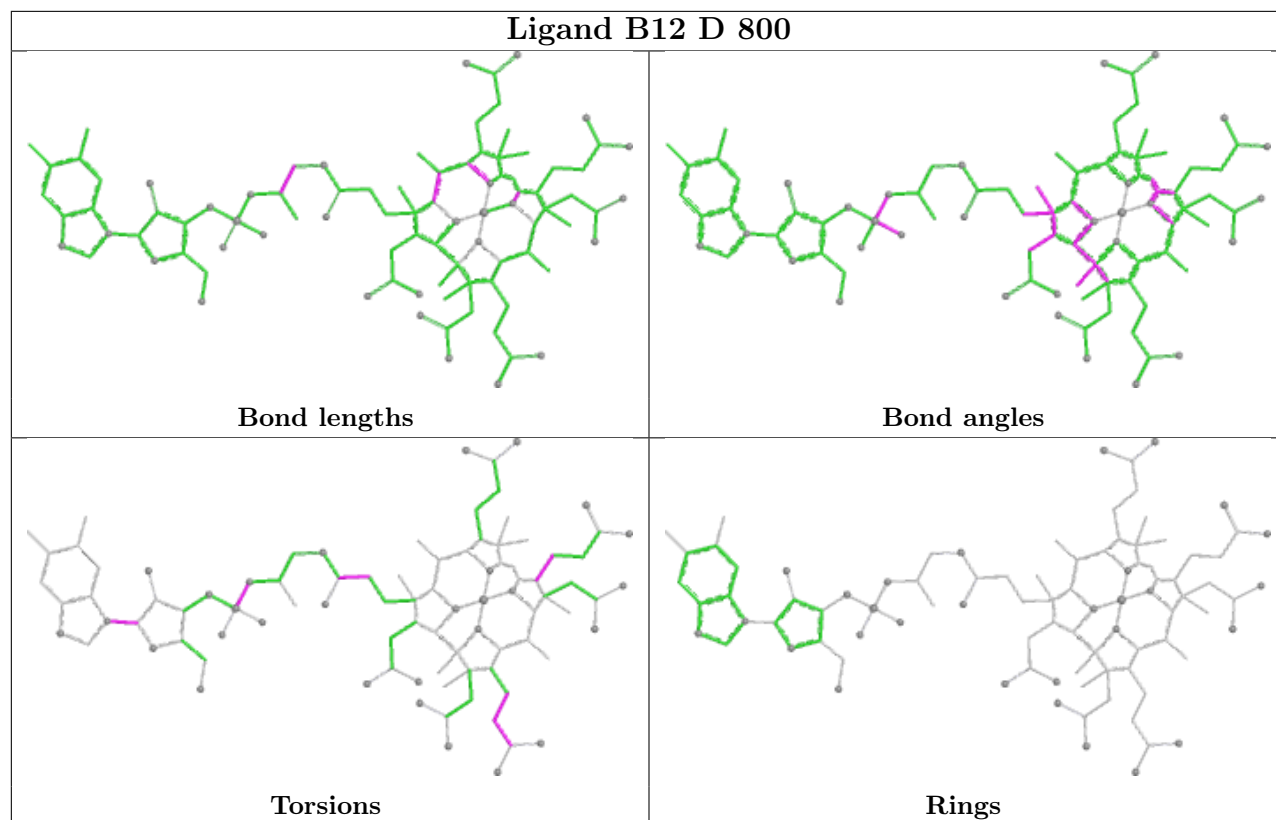


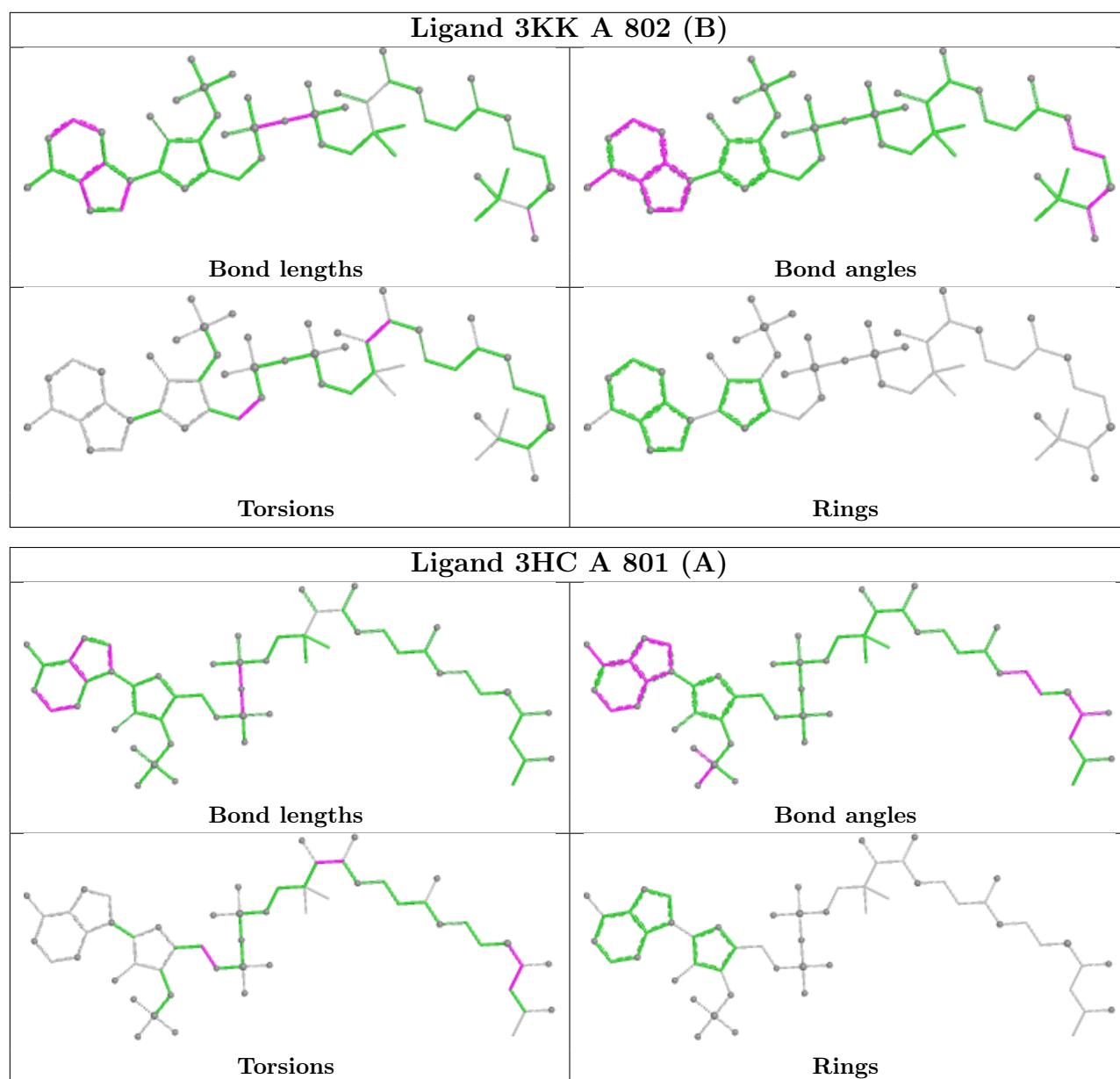
Ligand 5AD A 803 (C)



Ligand 3KK B 802 (B)







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(Å²)	Q<0.9	
1	A	557/584 (95%)	-0.75	0	100	100	16, 34, 62, 97	0
1	B	555/584 (95%)	-0.20	0	100	100	42, 56, 81, 101	0
2	C	131/158 (82%)	-0.69	0	100	100	19, 36, 61, 73	0
2	D	129/158 (81%)	0.35	3 (2%)	61	57	54, 77, 97, 105	0
All	All	1372/1484 (92%)	-0.42	3 (0%)	91	89	16, 49, 82, 105	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	132	ALA	2.5
2	D	4	THR	2.3
2	D	67	GLY	2.2

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

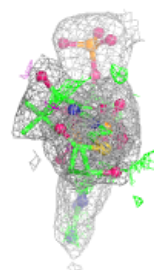
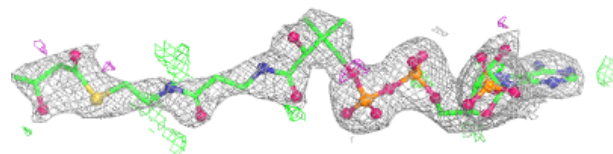
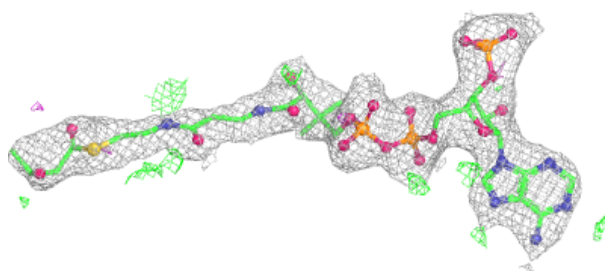
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	3HC	B	801[A]	54/54	0.92	0.10	45,55,63,65	54
4	3KK	B	802[B]	54/54	0.92	0.10	45,56,63,65	54
5	5AD	A	803[C]	18/18	0.92	0.15	17,27,31,32	18
6	B12	D	800	91/91	0.95	0.08	29,44,66,70	0
4	3KK	A	802[B]	54/54	0.96	0.07	6,32,45,47	54
3	3HC	A	801[A]	54/54	0.97	0.07	20,33,45,47	54
6	B12	C	800	91/91	0.98	0.06	14,24,35,38	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.

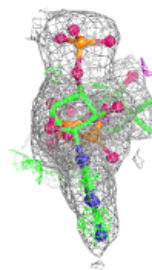
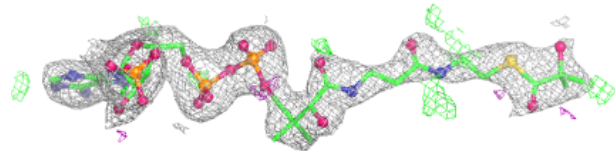
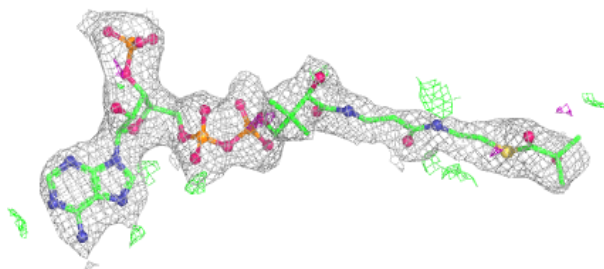
Electron density around 3HC B 801 (A):

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

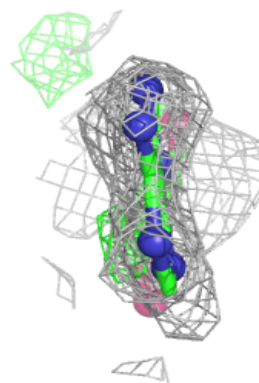
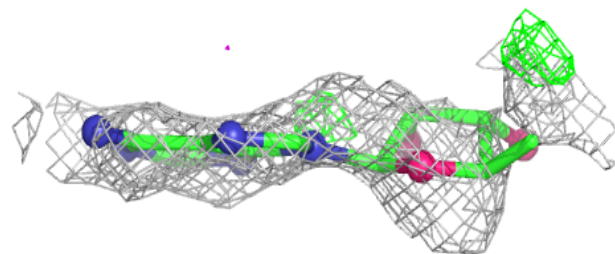
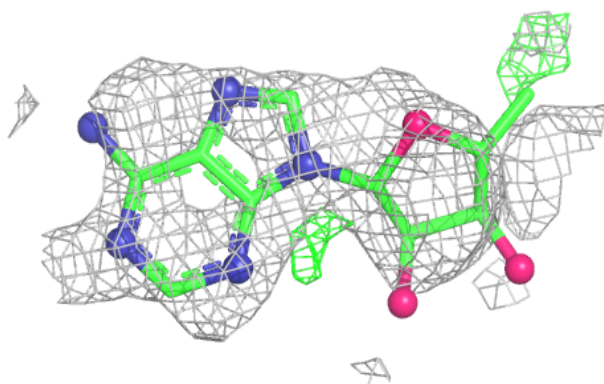


Electron density around 3KK B 802 (B):

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

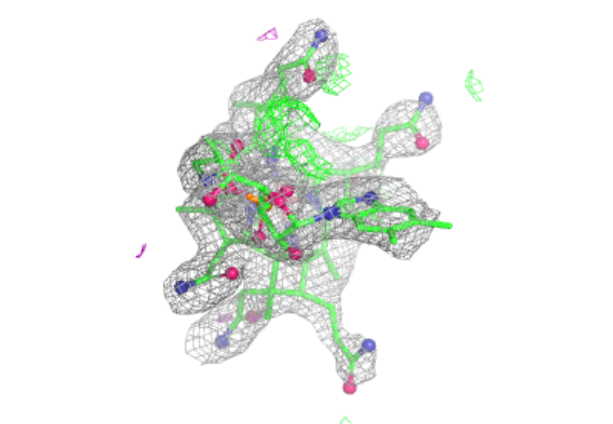
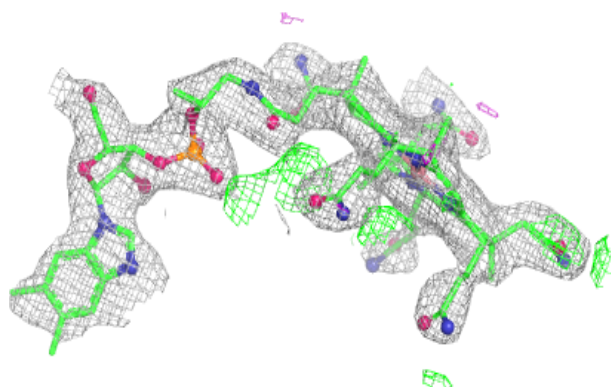
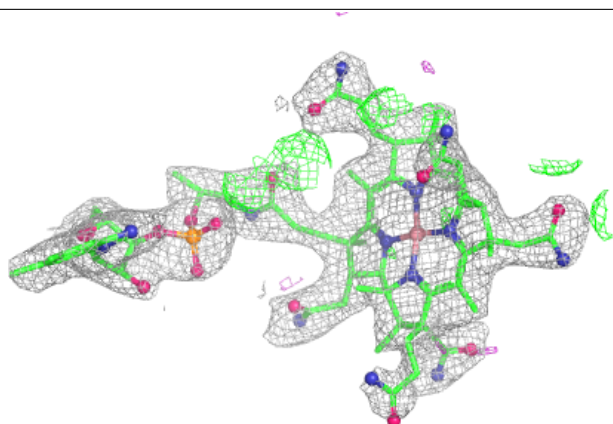
**Electron density around 5AD A 803 (C):**

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

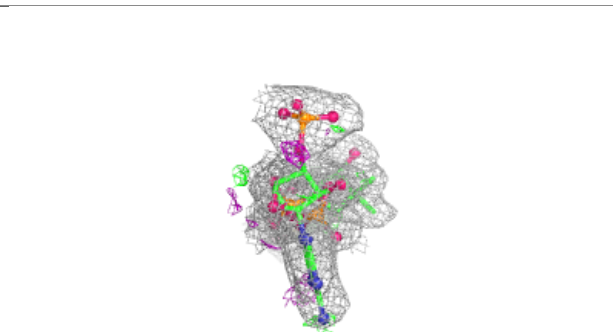
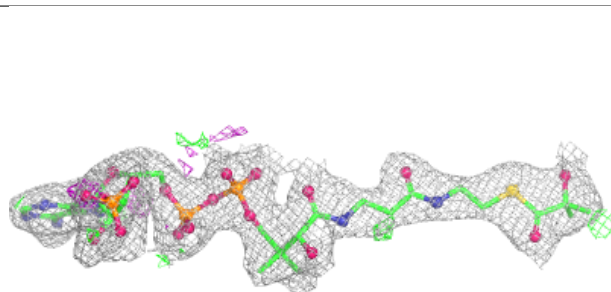
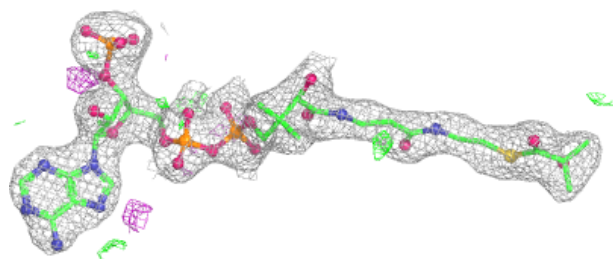


Electron density around B12 D 800:

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

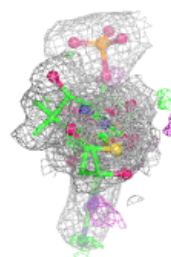
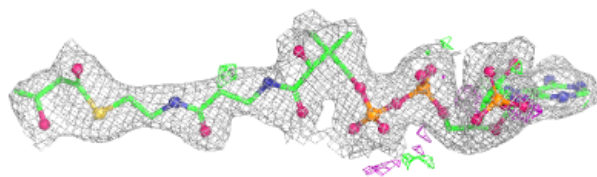
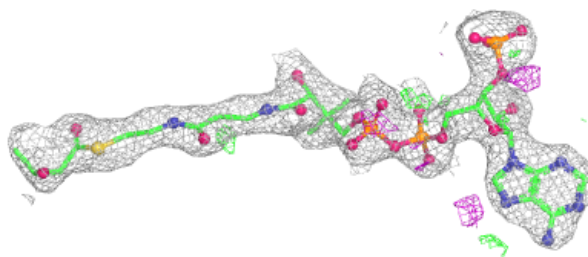
**Electron density around 3KK A 802 (B):**

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



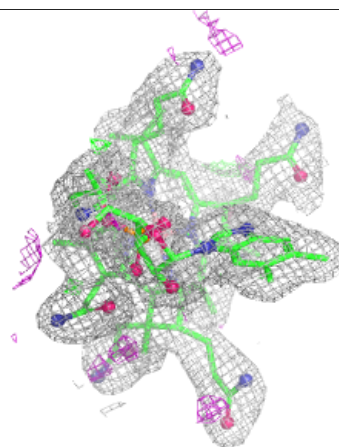
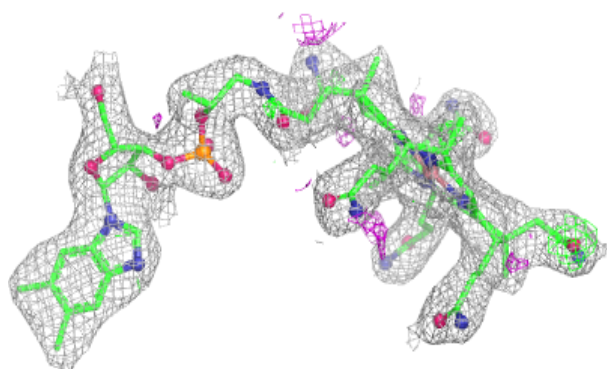
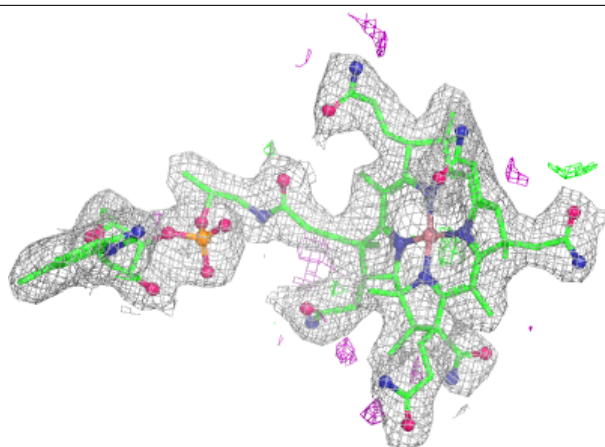
Electron density around 3HC A 801 (A):

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



Electron density around B12 C 800:

$2mF_o-DF_c$ (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.