



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:32 AM UTC

PDB ID : 5D6S / pdb_00005d6s
Title : Structure of epoxyqueuosine reductase from *Streptococcus thermophilus*.
Authors : Payne, K.A.P.; Fisher, K.; Dunstan, M.S.; Sjuts, H.; Leys, D.
Deposited on : 2015-08-12
Resolution : 2.65 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

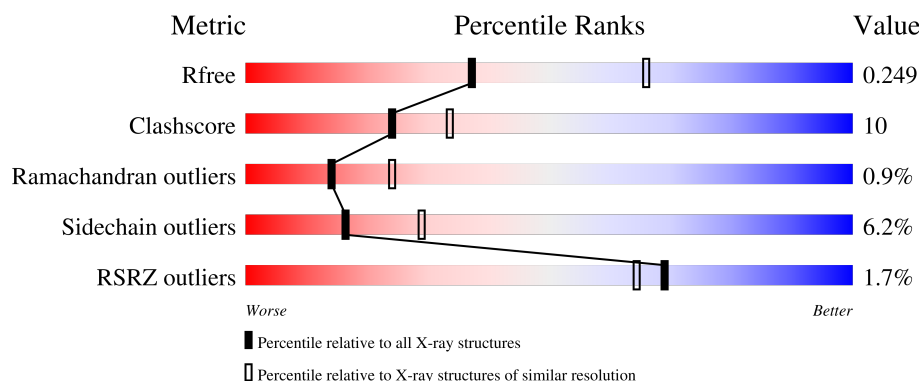
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	399	0.2% 75% 14% 7%
1	B	399	0.2% 77% 12% 7%
1	C	399	4% 76% 12% 8%
1	D	399	0.2% 76% 13% 7%
1	E	399	2% 76% 14% 7%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SF4	D	402	-	-	X	-
3	B12	A	403	-	-	X	-
3	B12	B	403	-	-	X	-
3	B12	E	403	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15145 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 Epoxyqueuosine reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	370	Total	C	N	O	S	0	0	0
			2912	1849	496	544	23			
1	B	370	Total	C	N	O	S	0	0	0
			2912	1849	496	544	23			
1	C	367	Total	C	N	O	S	0	0	0
			2888	1834	492	540	22			
1	D	370	Total	C	N	O	S	0	0	0
			2912	1849	496	544	23			
1	E	370	Total	C	N	O	S	0	0	0
			2912	1849	496	544	23			

There are 145 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-26	MET	-	initiating methionine	UNP A0A0F7K4Z9
A	-25	VAL	-	expression tag	UNP A0A0F7K4Z9
A	-24	GLN	-	expression tag	UNP A0A0F7K4Z9
A	-23	THR	-	expression tag	UNP A0A0F7K4Z9
A	-22	SER	-	expression tag	UNP A0A0F7K4Z9
A	-21	PHE	-	expression tag	UNP A0A0F7K4Z9
A	-20	GLU	-	expression tag	UNP A0A0F7K4Z9
A	-19	HIS	-	expression tag	UNP A0A0F7K4Z9
A	-18	HIS	-	expression tag	UNP A0A0F7K4Z9
A	-17	HIS	-	expression tag	UNP A0A0F7K4Z9
A	-16	HIS	-	expression tag	UNP A0A0F7K4Z9
A	-15	HIS	-	expression tag	UNP A0A0F7K4Z9
A	-14	HIS	-	expression tag	UNP A0A0F7K4Z9
A	-13	SER	-	expression tag	UNP A0A0F7K4Z9
A	-12	ALA	-	expression tag	UNP A0A0F7K4Z9
A	-11	GLY	-	expression tag	UNP A0A0F7K4Z9
A	-10	GLU	-	expression tag	UNP A0A0F7K4Z9
A	-9	ASN	-	expression tag	UNP A0A0F7K4Z9
A	-8	LEU	-	expression tag	UNP A0A0F7K4Z9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	TYR	-	expression tag	UNP A0A0F7K4Z9
A	-6	PHE	-	expression tag	UNP A0A0F7K4Z9
A	-5	GLN	-	expression tag	UNP A0A0F7K4Z9
A	-4	GLY	-	expression tag	UNP A0A0F7K4Z9
A	-3	ALA	-	expression tag	UNP A0A0F7K4Z9
A	-2	GLN	-	expression tag	UNP A0A0F7K4Z9
A	-1	ILE	-	expression tag	UNP A0A0F7K4Z9
A	0	SER	-	expression tag	UNP A0A0F7K4Z9
A	258	ASP	GLU	conflict	UNP A0A0F7K4Z9
A	342	GLY	ALA	conflict	UNP A0A0F7K4Z9
B	-26	MET	-	initiating methionine	UNP A0A0F7K4Z9
B	-25	VAL	-	expression tag	UNP A0A0F7K4Z9
B	-24	GLN	-	expression tag	UNP A0A0F7K4Z9
B	-23	THR	-	expression tag	UNP A0A0F7K4Z9
B	-22	SER	-	expression tag	UNP A0A0F7K4Z9
B	-21	PHE	-	expression tag	UNP A0A0F7K4Z9
B	-20	GLU	-	expression tag	UNP A0A0F7K4Z9
B	-19	HIS	-	expression tag	UNP A0A0F7K4Z9
B	-18	HIS	-	expression tag	UNP A0A0F7K4Z9
B	-17	HIS	-	expression tag	UNP A0A0F7K4Z9
B	-16	HIS	-	expression tag	UNP A0A0F7K4Z9
B	-15	HIS	-	expression tag	UNP A0A0F7K4Z9
B	-14	HIS	-	expression tag	UNP A0A0F7K4Z9
B	-13	SER	-	expression tag	UNP A0A0F7K4Z9
B	-12	ALA	-	expression tag	UNP A0A0F7K4Z9
B	-11	GLY	-	expression tag	UNP A0A0F7K4Z9
B	-10	GLU	-	expression tag	UNP A0A0F7K4Z9
B	-9	ASN	-	expression tag	UNP A0A0F7K4Z9
B	-8	LEU	-	expression tag	UNP A0A0F7K4Z9
B	-7	TYR	-	expression tag	UNP A0A0F7K4Z9
B	-6	PHE	-	expression tag	UNP A0A0F7K4Z9
B	-5	GLN	-	expression tag	UNP A0A0F7K4Z9
B	-4	GLY	-	expression tag	UNP A0A0F7K4Z9
B	-3	ALA	-	expression tag	UNP A0A0F7K4Z9
B	-2	GLN	-	expression tag	UNP A0A0F7K4Z9
B	-1	ILE	-	expression tag	UNP A0A0F7K4Z9
B	0	SER	-	expression tag	UNP A0A0F7K4Z9
B	258	ASP	GLU	conflict	UNP A0A0F7K4Z9
B	342	GLY	ALA	conflict	UNP A0A0F7K4Z9
C	-26	MET	-	initiating methionine	UNP A0A0F7K4Z9
C	-25	VAL	-	expression tag	UNP A0A0F7K4Z9
C	-24	GLN	-	expression tag	UNP A0A0F7K4Z9

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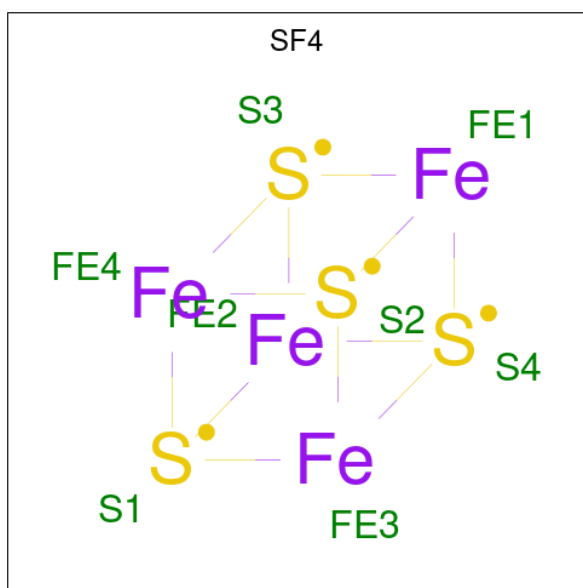
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C	-23	THR	-	expression tag	UNP A0A0F7K4Z9
C	-22	SER	-	expression tag	UNP A0A0F7K4Z9
C	-21	PHE	-	expression tag	UNP A0A0F7K4Z9
C	-20	GLU	-	expression tag	UNP A0A0F7K4Z9
C	-19	HIS	-	expression tag	UNP A0A0F7K4Z9
C	-18	HIS	-	expression tag	UNP A0A0F7K4Z9
C	-17	HIS	-	expression tag	UNP A0A0F7K4Z9
C	-16	HIS	-	expression tag	UNP A0A0F7K4Z9
C	-15	HIS	-	expression tag	UNP A0A0F7K4Z9
C	-14	HIS	-	expression tag	UNP A0A0F7K4Z9
C	-13	SER	-	expression tag	UNP A0A0F7K4Z9
C	-12	ALA	-	expression tag	UNP A0A0F7K4Z9
C	-11	GLY	-	expression tag	UNP A0A0F7K4Z9
C	-10	GLU	-	expression tag	UNP A0A0F7K4Z9
C	-9	ASN	-	expression tag	UNP A0A0F7K4Z9
C	-8	LEU	-	expression tag	UNP A0A0F7K4Z9
C	-7	TYR	-	expression tag	UNP A0A0F7K4Z9
C	-6	PHE	-	expression tag	UNP A0A0F7K4Z9
C	-5	GLN	-	expression tag	UNP A0A0F7K4Z9
C	-4	GLY	-	expression tag	UNP A0A0F7K4Z9
C	-3	ALA	-	expression tag	UNP A0A0F7K4Z9
C	-2	GLN	-	expression tag	UNP A0A0F7K4Z9
C	-1	ILE	-	expression tag	UNP A0A0F7K4Z9
C	0	SER	-	expression tag	UNP A0A0F7K4Z9
C	260	ASP	GLU	conflict	UNP A0A0F7K4Z9
C	344	GLY	ALA	conflict	UNP A0A0F7K4Z9
D	-26	MET	-	initiating methionine	UNP A0A0F7K4Z9
D	-25	VAL	-	expression tag	UNP A0A0F7K4Z9
D	-24	GLN	-	expression tag	UNP A0A0F7K4Z9
D	-23	THR	-	expression tag	UNP A0A0F7K4Z9
D	-22	SER	-	expression tag	UNP A0A0F7K4Z9
D	-21	PHE	-	expression tag	UNP A0A0F7K4Z9
D	-20	GLU	-	expression tag	UNP A0A0F7K4Z9
D	-19	HIS	-	expression tag	UNP A0A0F7K4Z9
D	-18	HIS	-	expression tag	UNP A0A0F7K4Z9
D	-17	HIS	-	expression tag	UNP A0A0F7K4Z9
D	-16	HIS	-	expression tag	UNP A0A0F7K4Z9
D	-15	HIS	-	expression tag	UNP A0A0F7K4Z9
D	-14	HIS	-	expression tag	UNP A0A0F7K4Z9
D	-13	SER	-	expression tag	UNP A0A0F7K4Z9
D	-12	ALA	-	expression tag	UNP A0A0F7K4Z9
D	-11	GLY	-	expression tag	UNP A0A0F7K4Z9

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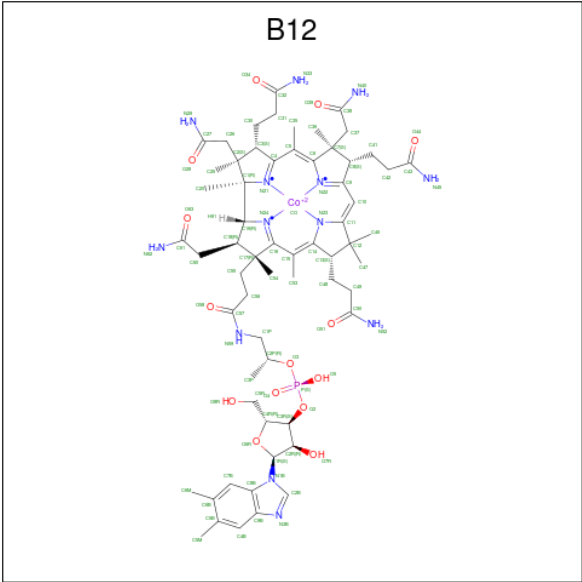
Chain	Residue	Modelled	Actual	Comment	Reference
D	-10	GLU	-	expression tag	UNP A0A0F7K4Z9
D	-9	ASN	-	expression tag	UNP A0A0F7K4Z9
D	-8	LEU	-	expression tag	UNP A0A0F7K4Z9
D	-7	TYR	-	expression tag	UNP A0A0F7K4Z9
D	-6	PHE	-	expression tag	UNP A0A0F7K4Z9
D	-5	GLN	-	expression tag	UNP A0A0F7K4Z9
D	-4	GLY	-	expression tag	UNP A0A0F7K4Z9
D	-3	ALA	-	expression tag	UNP A0A0F7K4Z9
D	-2	GLN	-	expression tag	UNP A0A0F7K4Z9
D	-1	ILE	-	expression tag	UNP A0A0F7K4Z9
D	0	SER	-	expression tag	UNP A0A0F7K4Z9
D	258	ASP	GLU	conflict	UNP A0A0F7K4Z9
D	342	GLY	ALA	conflict	UNP A0A0F7K4Z9
E	-26	MET	-	initiating methionine	UNP A0A0F7K4Z9
E	-25	VAL	-	expression tag	UNP A0A0F7K4Z9
E	-24	GLN	-	expression tag	UNP A0A0F7K4Z9
E	-23	THR	-	expression tag	UNP A0A0F7K4Z9
E	-22	SER	-	expression tag	UNP A0A0F7K4Z9
E	-21	PHE	-	expression tag	UNP A0A0F7K4Z9
E	-20	GLU	-	expression tag	UNP A0A0F7K4Z9
E	-19	HIS	-	expression tag	UNP A0A0F7K4Z9
E	-18	HIS	-	expression tag	UNP A0A0F7K4Z9
E	-17	HIS	-	expression tag	UNP A0A0F7K4Z9
E	-16	HIS	-	expression tag	UNP A0A0F7K4Z9
E	-15	HIS	-	expression tag	UNP A0A0F7K4Z9
E	-14	HIS	-	expression tag	UNP A0A0F7K4Z9
E	-13	SER	-	expression tag	UNP A0A0F7K4Z9
E	-12	ALA	-	expression tag	UNP A0A0F7K4Z9
E	-11	GLY	-	expression tag	UNP A0A0F7K4Z9
E	-10	GLU	-	expression tag	UNP A0A0F7K4Z9
E	-9	ASN	-	expression tag	UNP A0A0F7K4Z9
E	-8	LEU	-	expression tag	UNP A0A0F7K4Z9
E	-7	TYR	-	expression tag	UNP A0A0F7K4Z9
E	-6	PHE	-	expression tag	UNP A0A0F7K4Z9
E	-5	GLN	-	expression tag	UNP A0A0F7K4Z9
E	-4	GLY	-	expression tag	UNP A0A0F7K4Z9
E	-3	ALA	-	expression tag	UNP A0A0F7K4Z9
E	-2	GLN	-	expression tag	UNP A0A0F7K4Z9
E	-1	ILE	-	expression tag	UNP A0A0F7K4Z9
E	0	SER	-	expression tag	UNP A0A0F7K4Z9
E	258	ASP	GLU	conflict	UNP A0A0F7K4Z9
E	342	GLY	ALA	conflict	UNP A0A0F7K4Z9

- Molecule 2 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			8	4	4		
2	A	1	Total	Fe	S	0	0
			8	4	4		
2	B	1	Total	Fe	S	0	0
			8	4	4		
2	B	1	Total	Fe	S	0	0
			8	4	4		
2	C	1	Total	Fe	S	0	0
			8	4	4		
2	C	1	Total	Fe	S	0	0
			8	4	4		
2	D	1	Total	Fe	S	0	0
			8	4	4		
2	D	1	Total	Fe	S	0	0
			8	4	4		
2	E	1	Total	Fe	S	0	0
			8	4	4		
2	E	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 3 is COBALAMIN (CCD ID: B12) (formula: $\text{C}_{62}\text{H}_{89}\text{CoN}_{13}\text{O}_{14}\text{P}$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total 91	C 62	Co 1	N 13	O 14	P 1	0	0
3	B	1	Total 91	C 62	Co 1	N 13	O 14	P 1	0	0
3	C	1	Total 91	C 62	Co 1	N 13	O 14	P 1	0	0
3	D	1	Total 91	C 62	Co 1	N 13	O 14	P 1	0	0
3	E	1	Total 91	C 62	Co 1	N 13	O 14	P 1	0	0

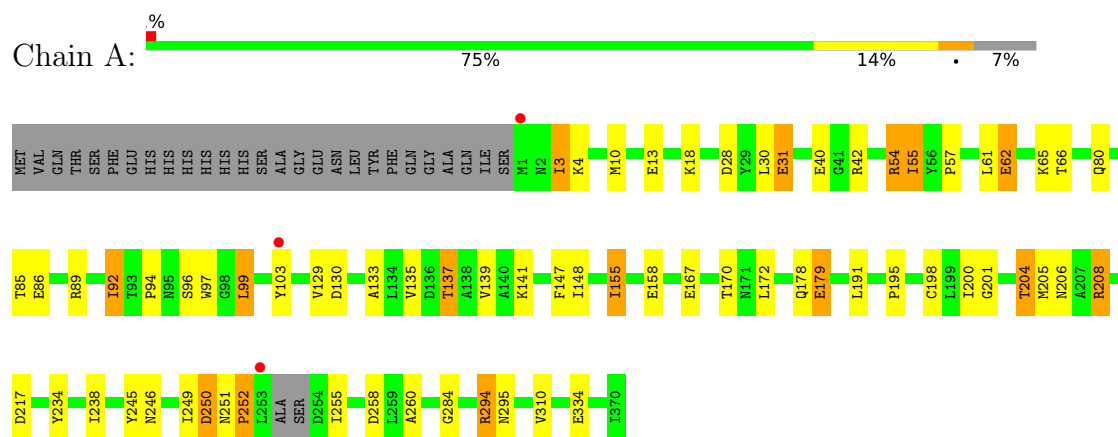
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	19	Total	O	0	0
			19	19		
4	B	14	Total	O	0	0
			14	14		
4	C	15	Total	O	0	0
			15	15		
4	D	14	Total	O	0	0
			14	14		
4	E	12	Total	O	0	0
			12	12		

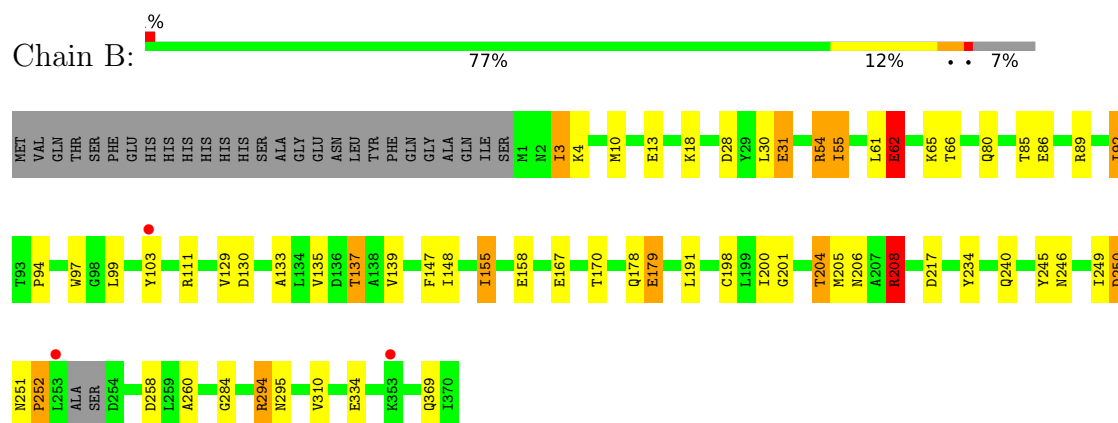
3 Residue-property plots

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.

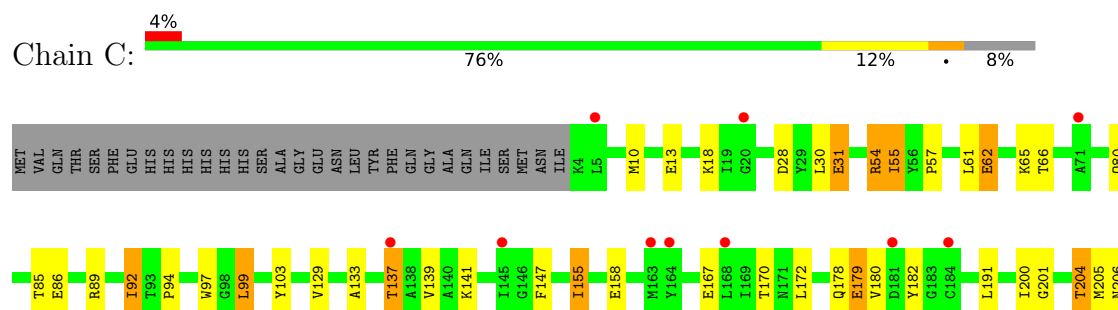
• Molecule 1: Epoxyqueuosine reductase

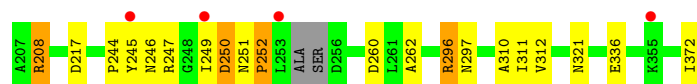


• Molecule 1: Epoxyqueuosine reductase

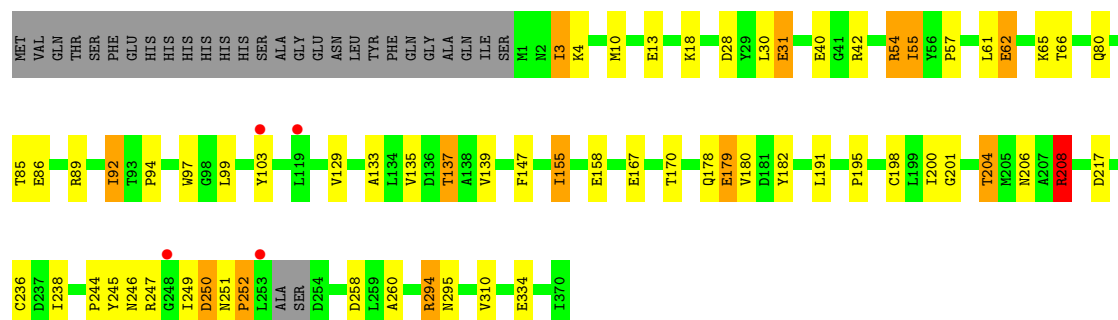
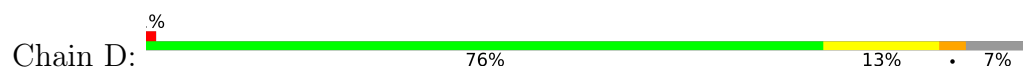


• Molecule 1: Epoxyqueuosine reductase

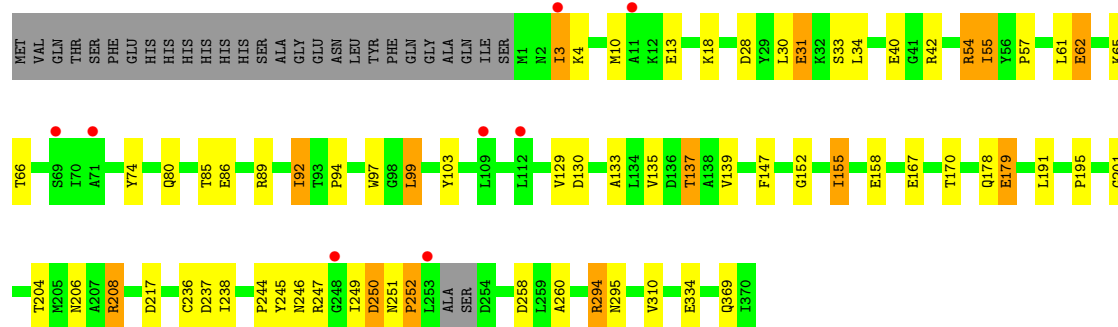
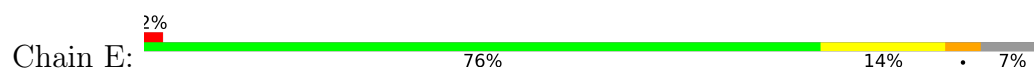




• Molecule 1: Epoxyqueuosine reductase



• Molecule 1: Epoxyqueuosine reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	106.12Å 106.12Å 332.28Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	88.58 – 2.65 88.58 – 2.65	Depositor EDS
% Data completeness (in resolution range)	98.8 (88.58-2.65) 98.8 (88.58-2.65)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 2.65Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.214 , 0.253 0.214 , 0.249	Depositor DCC
R_{free} test set	3163 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	65.9	Xtriage
Anisotropy	0.082	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 36.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15145	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.56% 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: B12, SF4

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	0.73	0/2965	0.94	0/3993
1	B	0.76	1/2965 (0.0%)	0.94	0/3993
1	C	0.73	0/2941	0.93	0/3961
1	D	0.72	0/2965	0.92	0/3993
1	E	0.68	0/2965	0.91	0/3993
All	All	0.72	1/14801 (0.0%)	0.93	0/19933

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	62	GLU	N-CA	5.09	1.52	1.46

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	2912	0	2929	51	0
1	B	2912	0	2929	46	0
1	C	2888	0	2900	44	0
1	D	2912	0	2929	45	0
1	E	2912	0	2929	53	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	16	0	0	1	0
2	B	16	0	0	1	0
2	C	16	0	0	0	0
2	D	16	0	0	2	0
2	E	16	0	0	1	0
3	A	91	0	88	24	0
3	B	91	0	88	22	0
3	C	91	0	88	16	0
3	D	91	0	88	16	0
3	E	91	0	88	22	0
4	A	19	0	0	0	0
4	B	14	0	0	1	0
4	C	15	0	0	0	0
4	D	14	0	0	0	0
4	E	12	0	0	3	0
All	All	15145	0	15056	300	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 10.

All (300) 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:148:ILE:CD1	3:B:403:B12:H2B	1.78	1.13
1:B:148:ILE:HD13	3:B:403:B12:H2B	1.31	1.12
1:A:148:ILE:HD12	3:A:403:B12:N3B	1.79	0.96
3:C:403:B12:H362	3:C:403:B12:H351	1.52	0.92
1:E:34:LEU:HD23	3:E:403:B12:H3P1	1.52	0.89
1:D:54:ARG:NH1	3:D:403:B12:O58	2.05	0.87
1:C:137:THR:HG21	3:C:403:B12:O8R	1.74	0.87
1:B:54:ARG:NH1	3:B:403:B12:O58	2.08	0.85
3:D:403:B12:H362	3:D:403:B12:H351	1.58	0.83
3:D:403:B12:H552	3:D:403:B12:H531	1.60	0.82
1:E:34:LEU:CD2	3:E:403:B12:C3P	2.56	0.82
1:B:148:ILE:HD12	3:B:403:B12:H2B	1.60	0.81
1:E:34:LEU:CD2	3:E:403:B12:H3P1	2.11	0.80
1:E:54:ARG:NH1	3:E:403:B12:O58	2.11	0.80
1:C:141:LYS:NZ	3:C:403:B12:H4B	1.96	0.80
1:C:205:MET:SD	3:C:403:B12:N52	2.54	0.79
1:A:148:ILE:CD1	3:A:403:B12:N3B	2.46	0.79
3:B:403:B12:H362	3:B:403:B12:H351	1.62	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:74:TYR:OH	4:E:501:HOH:O	1.99	0.79
1:A:148:ILE:HD13	3:A:403:B12:C2B	2.12	0.78
3:A:403:B12:H351	3:A:403:B12:H362	1.66	0.77
1:C:141:LYS:HZ3	3:C:403:B12:H4B	1.48	0.77
1:A:148:ILE:CD1	3:A:403:B12:C2B	2.63	0.77
3:E:403:B12:H362	3:E:403:B12:H351	1.67	0.76
1:A:137:THR:HG21	3:A:403:B12:O8R	1.87	0.74
1:B:294:ARG:NH1	1:B:295:ASN:OD1	2.20	0.73
1:D:294:ARG:NH1	1:D:295:ASN:OD1	2.21	0.73
3:C:403:B12:O63	3:C:403:B12:H551	1.89	0.73
1:A:141:LYS:NZ	3:A:403:B12:H4B	2.04	0.72
1:E:294:ARG:NH1	1:E:295:ASN:OD1	2.22	0.72
1:A:294:ARG:NH1	1:A:295:ASN:OD1	2.23	0.72
1:E:34:LEU:HD21	3:E:403:B12:H3P3	1.72	0.71
1:C:296:ARG:NH1	1:C:297:ASN:OD1	2.24	0.71
3:E:403:B12:H3P2	3:E:403:B12:O2	1.92	0.70
3:C:403:B12:H421	3:C:403:B12:C10	2.21	0.70
1:B:137:THR:HG21	3:B:403:B12:O8R	1.91	0.69
1:E:34:LEU:CD2	3:E:403:B12:H3P3	2.20	0.69
1:E:130:ASP:O	3:E:403:B12:N62	2.26	0.68
3:C:403:B12:H473	3:C:403:B12:H491	1.74	0.68
3:B:403:B12:H421	3:B:403:B12:C10	2.23	0.68
1:E:294:ARG:HG2	1:E:294:ARG:HH11	1.57	0.67
1:A:294:ARG:HH11	1:A:294:ARG:HG2	1.58	0.67
3:D:403:B12:H531	3:D:403:B12:C55	2.22	0.66
1:A:130:ASP:OD1	3:A:403:B12:H542	1.96	0.66
1:D:294:ARG:HG2	1:D:294:ARG:HH11	1.59	0.66
1:C:296:ARG:HG2	1:C:296:ARG:HH11	1.60	0.65
1:B:294:ARG:HG2	1:B:294:ARG:HH11	1.62	0.64
1:E:33:SER:OG	4:E:502:HOH:O	2.15	0.64
1:A:141:LYS:HZ2	3:A:403:B12:H4B	1.61	0.64
1:E:34:LEU:HD21	3:E:403:B12:C3P	2.29	0.63
3:C:403:B12:H362	3:C:403:B12:C35	2.26	0.63
1:C:372:ILE:O	1:C:372:ILE:HG22	1.97	0.62
1:A:206:ASN:C	1:A:206:ASN:HD22	2.07	0.61
3:A:403:B12:C10	3:A:403:B12:H421	2.30	0.61
1:B:85:THR:HG22	1:B:86:GLU:H	1.69	0.58
1:D:206:ASN:C	1:D:206:ASN:HD22	2.12	0.58
1:B:246:ASN:HA	1:B:249:ILE:CD1	2.33	0.58
1:B:206:ASN:C	1:B:206:ASN:HD22	2.11	0.58
1:A:246:ASN:HA	1:A:249:ILE:CD1	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:246:ASN:HA	1:D:249:ILE:CD1	2.33	0.57
1:C:246:ASN:HA	1:C:249:ILE:CD1	2.35	0.57
1:C:85:THR:HG22	1:C:86:GLU:H	1.69	0.57
3:A:403:B12:H311	3:A:403:B12:H353	1.87	0.57
1:B:198:CYS:SG	2:B:402:SF4:S2	3.02	0.57
3:E:403:B12:H362	3:E:403:B12:C35	2.34	0.57
1:C:61:LEU:O	1:C:62:GLU:CB	2.52	0.57
3:B:403:B12:H533	3:B:403:B12:H481	1.87	0.56
3:D:403:B12:C10	3:D:403:B12:H421	2.34	0.56
1:E:85:THR:HG22	1:E:86:GLU:H	1.70	0.56
1:E:246:ASN:HA	1:E:249:ILE:CD1	2.34	0.56
1:D:85:THR:HG22	1:D:86:GLU:H	1.69	0.56
3:C:403:B12:H551	3:C:403:B12:C61	2.36	0.56
1:C:147:PHE:CE1	1:C:155:ILE:HB	2.41	0.55
3:C:403:B12:H492	3:C:403:B12:N23	2.21	0.55
3:A:403:B12:H362	3:A:403:B12:C35	2.36	0.55
1:C:217:ASP:O	1:C:296:ARG:NH2	2.39	0.55
3:D:403:B12:H481	3:D:403:B12:H533	1.87	0.55
3:D:403:B12:H491	3:D:403:B12:H473	1.87	0.55
1:E:152:GLY:O	3:E:403:B12:O8R	2.25	0.55
1:A:85:THR:HG22	1:A:86:GLU:H	1.72	0.55
3:A:403:B12:H531	3:A:403:B12:H552	1.89	0.54
1:A:217:ASP:O	1:A:294:ARG:NH2	2.41	0.54
3:A:403:B12:H531	3:A:403:B12:C55	2.36	0.54
1:E:206:ASN:C	1:E:206:ASN:HD22	2.16	0.54
1:E:217:ASP:O	1:E:294:ARG:NH2	2.39	0.54
1:B:217:ASP:O	1:B:294:ARG:NH2	2.40	0.54
1:B:130:ASP:OD1	3:B:403:B12:H542	2.08	0.53
3:B:403:B12:H252	3:B:403:B12:H601	1.90	0.53
3:B:403:B12:H473	3:B:403:B12:H491	1.90	0.53
1:A:147:PHE:CE1	1:A:155:ILE:HB	2.43	0.53
3:A:403:B12:H3	3:A:403:B12:O28	2.08	0.53
1:B:147:PHE:CE1	1:B:155:ILE:HB	2.43	0.53
3:E:403:B12:N62	3:E:403:B12:H541	2.23	0.53
1:D:147:PHE:CE1	1:D:155:ILE:HB	2.45	0.53
1:E:147:PHE:CE1	1:E:155:ILE:HB	2.44	0.53
1:B:139:VAL:HG11	1:B:167:GLU:HG2	1.91	0.52
1:D:206:ASN:HD21	1:D:208:ARG:HB2	1.74	0.52
1:D:250:ASP:C	1:D:250:ASP:OD1	2.52	0.52
3:A:403:B12:H473	3:A:403:B12:H491	1.92	0.52
3:A:403:B12:H561	3:A:403:B12:C53	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:250:ASP:C	1:B:250:ASP:OD1	2.52	0.52
1:C:139:VAL:HG11	1:C:167:GLU:HG2	1.92	0.52
1:D:137:THR:HG21	3:D:403:B12:O8R	2.09	0.52
1:D:217:ASP:O	1:D:294:ARG:NH2	2.42	0.52
1:D:206:ASN:ND2	1:D:208:ARG:HB2	2.24	0.52
1:E:250:ASP:OD1	1:E:250:ASP:C	2.52	0.52
1:A:250:ASP:OD1	1:A:250:ASP:C	2.53	0.51
1:A:139:VAL:HG11	1:A:167:GLU:HG2	1.92	0.51
1:D:139:VAL:HG11	1:D:167:GLU:HG2	1.92	0.51
1:E:61:LEU:O	1:E:62:GLU:CB	2.59	0.51
1:E:10:MET:HA	1:E:13:GLU:HB3	1.93	0.51
1:E:139:VAL:HG11	1:E:167:GLU:HG2	1.93	0.51
1:D:10:MET:HA	1:D:13:GLU:HB3	1.93	0.51
1:C:10:MET:HA	1:C:13:GLU:HB3	1.93	0.50
3:B:403:B12:H2R	3:B:403:B12:H7B	1.91	0.50
3:D:403:B12:H362	3:D:403:B12:C35	2.35	0.50
3:E:403:B12:C3P	3:E:403:B12:O2	2.59	0.50
1:A:158:GLU:OE1	1:A:158:GLU:N	2.43	0.50
1:E:31:GLU:CG	1:E:55:ILE:HD13	2.41	0.50
1:A:61:LEU:O	1:A:62:GLU:CB	2.59	0.50
1:C:206:ASN:HD22	1:C:206:ASN:C	2.18	0.50
1:C:250:ASP:OD1	1:C:250:ASP:C	2.54	0.50
1:A:206:ASN:ND2	1:A:208:ARG:HB2	2.26	0.50
1:B:206:ASN:ND2	1:B:208:ARG:HB2	2.26	0.50
1:B:31:GLU:CG	1:B:55:ILE:HD13	2.42	0.49
1:C:206:ASN:ND2	1:C:208:ARG:HB2	2.27	0.49
1:B:245:TYR:O	1:B:249:ILE:HD11	2.12	0.49
1:C:89:ARG:NH1	1:C:260:ASP:OD1	2.45	0.49
1:C:245:TYR:O	1:C:249:ILE:HD11	2.12	0.49
1:D:31:GLU:CG	1:D:55:ILE:HD13	2.43	0.49
1:E:245:TYR:O	1:E:249:ILE:HD11	2.13	0.49
1:D:89:ARG:NH1	1:D:258:ASP:OD1	2.46	0.49
1:E:236:CYS:HB3	2:E:402:SF4:S4	2.52	0.49
1:C:206:ASN:HD21	1:C:208:ARG:HB2	1.77	0.49
1:D:245:TYR:O	1:D:249:ILE:HD11	2.13	0.49
1:A:245:TYR:O	1:A:249:ILE:HD11	2.12	0.49
1:B:10:MET:HA	1:B:13:GLU:HB3	1.94	0.49
1:C:31:GLU:CG	1:C:55:ILE:HD13	2.42	0.48
1:A:31:GLU:CG	1:A:55:ILE:HD13	2.43	0.48
1:A:206:ASN:HD21	1:A:208:ARG:HB2	1.78	0.48
3:B:403:B12:H362	3:B:403:B12:C35	2.37	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:85:THR:HG22	1:C:86:GLU:N	2.28	0.48
1:A:3:ILE:HD12	1:A:4:LYS:H	1.78	0.48
1:B:66:THR:HG23	1:B:170:THR:HG23	1.95	0.48
1:B:85:THR:HG22	1:B:86:GLU:N	2.28	0.48
1:A:10:MET:HA	1:A:13:GLU:HB3	1.95	0.48
1:E:206:ASN:HD21	1:E:208:ARG:HB2	1.78	0.48
1:A:66:THR:HG23	1:A:170:THR:HG23	1.96	0.48
1:A:85:THR:HG22	1:A:86:GLU:N	2.28	0.48
3:A:403:B12:H481	3:A:403:B12:H533	1.96	0.47
1:E:85:THR:HG22	1:E:86:GLU:N	2.29	0.47
1:E:158:GLU:OE1	1:E:158:GLU:N	2.43	0.47
1:D:85:THR:HG22	1:D:86:GLU:N	2.29	0.47
1:D:236:CYS:HB3	2:D:402:SF4:S4	2.53	0.47
1:A:54:ARG:HH12	3:A:403:B12:C57	2.25	0.47
1:E:206:ASN:ND2	1:E:208:ARG:HB2	2.28	0.47
1:A:206:ASN:C	1:A:206:ASN:ND2	2.71	0.47
1:E:66:THR:HG23	1:E:170:THR:HG23	1.95	0.47
1:A:89:ARG:NH1	1:A:258:ASP:OD1	2.48	0.47
1:A:94:PRO:HA	1:A:97:TRP:CD2	2.50	0.47
1:C:296:ARG:HH11	1:C:296:ARG:CG	2.27	0.47
1:E:294:ARG:HH11	1:E:294:ARG:CG	2.26	0.47
3:E:403:B12:H421	3:E:403:B12:C10	2.45	0.47
1:B:206:ASN:HD21	1:B:208:ARG:HB2	1.79	0.47
1:C:158:GLU:OE1	1:C:158:GLU:N	2.42	0.47
1:E:89:ARG:NH1	1:E:258:ASP:OD1	2.48	0.47
3:E:403:B12:H262	3:E:403:B12:H91	1.77	0.47
1:B:31:GLU:HG2	1:B:55:ILE:HD13	1.98	0.46
1:B:89:ARG:NH1	1:B:258:ASP:OD1	2.49	0.46
1:D:66:THR:HG23	1:D:170:THR:HG23	1.96	0.46
1:A:198:CYS:SG	2:A:402:SF4:S2	3.13	0.46
1:B:158:GLU:OE1	1:B:158:GLU:N	2.45	0.46
3:C:403:B12:H531	3:C:403:B12:H552	1.98	0.46
1:C:66:THR:HG23	1:C:170:THR:HG23	1.96	0.46
1:E:3:ILE:HD12	1:E:4:LYS:H	1.79	0.46
3:E:403:B12:H312	3:E:403:B12:C25	2.46	0.46
1:B:61:LEU:O	1:B:62:GLU:CB	2.64	0.46
1:D:3:ILE:HD12	1:D:4:LYS:H	1.79	0.46
1:D:158:GLU:OE1	1:D:158:GLU:N	2.43	0.46
1:E:237:ASP:HA	4:E:510:HOH:O	2.16	0.46
3:D:403:B12:H8	3:D:403:B12:O39	2.15	0.46
1:E:31:GLU:HG2	1:E:55:ILE:HD13	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:LEU:O	1:A:31:GLU:C	2.59	0.46
1:B:3:ILE:HD12	1:B:4:LYS:H	1.81	0.46
1:D:94:PRO:HA	1:D:97:TRP:CD2	2.51	0.46
3:C:403:B12:O28	3:C:403:B12:H3	2.16	0.45
3:B:403:B12:H7B	3:B:403:B12:C2R	2.47	0.45
1:B:148:ILE:HD12	3:B:403:B12:C2B	2.38	0.45
1:C:251:ASN:HB3	1:C:252:PRO:CD	2.47	0.45
1:C:31:GLU:HG2	1:C:55:ILE:HD13	1.99	0.45
1:B:240:GLN:NE2	4:B:503:HOH:O	2.41	0.45
1:C:205:MET:HE1	3:C:403:B12:N45	2.31	0.45
3:E:403:B12:H312	3:E:403:B12:H251	1.98	0.45
1:D:294:ARG:HH11	1:D:294:ARG:CG	2.28	0.45
1:E:94:PRO:HA	1:E:97:TRP:CD2	2.52	0.45
1:B:30:LEU:O	1:B:31:GLU:C	2.60	0.45
1:D:251:ASN:HB3	1:D:252:PRO:CD	2.47	0.45
1:A:99:LEU:HD23	1:A:103:TYR:HD2	1.82	0.45
1:E:99:LEU:HD23	1:E:103:TYR:HD2	1.82	0.44
1:C:321:ASN:CG	1:C:321:ASN:O	2.60	0.44
3:A:403:B12:H253	3:A:403:B12:H301	1.84	0.44
3:A:403:B12:N23	3:A:403:B12:H492	2.33	0.44
1:D:206:ASN:C	1:D:206:ASN:ND2	2.75	0.44
1:D:31:GLU:HG2	1:D:55:ILE:HD13	2.00	0.44
1:E:251:ASN:HB3	1:E:252:PRO:CD	2.48	0.44
1:A:31:GLU:HG2	1:A:55:ILE:HD13	2.00	0.44
3:B:403:B12:H253	3:B:403:B12:H301	1.76	0.44
1:C:92:ILE:CG2	1:C:262:ALA:O	2.66	0.44
1:C:178:GLN:O	1:C:179:GLU:HB2	2.18	0.44
1:A:234:TYR:CZ	1:A:284:GLY:HA2	2.53	0.44
1:B:99:LEU:HD23	1:B:103:TYR:HD2	1.83	0.44
1:C:200:ILE:HB	1:C:204:THR:HG22	2.00	0.43
3:C:403:B12:H531	3:C:403:B12:C55	2.48	0.43
1:D:178:GLN:O	1:D:179:GLU:HB2	2.16	0.43
1:B:94:PRO:HA	1:B:97:TRP:CD2	2.53	0.43
1:C:191:LEU:HD21	1:C:201:GLY:HA2	2.00	0.43
1:D:30:LEU:O	1:D:31:GLU:C	2.62	0.43
1:B:92:ILE:CG2	1:B:260:ALA:O	2.66	0.43
1:B:130:ASP:HA	3:B:403:B12:H602	2.01	0.43
1:B:178:GLN:O	1:B:179:GLU:HB2	2.18	0.43
1:C:94:PRO:HA	1:C:97:TRP:CD2	2.54	0.43
1:D:61:LEU:O	1:D:62:GLU:CB	2.67	0.43
1:A:54:ARG:HD2	1:A:133:ALA:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:GLN:O	1:A:179:GLU:HB2	2.18	0.43
1:C:92:ILE:HD13	1:C:92:ILE:HG21	1.75	0.43
3:E:403:B12:H541	3:E:403:B12:C61	2.49	0.43
1:C:30:LEU:O	1:C:31:GLU:C	2.62	0.43
1:E:178:GLN:O	1:E:179:GLU:HB2	2.18	0.43
1:A:200:ILE:HB	1:A:204:THR:HG22	2.00	0.42
1:A:251:ASN:HB3	1:A:252:PRO:CD	2.49	0.42
1:B:251:ASN:HB3	1:B:252:PRO:CD	2.49	0.42
3:D:403:B12:C55	3:D:403:B12:C53	2.95	0.42
1:E:191:LEU:HD21	1:E:201:GLY:HA2	2.00	0.42
3:E:403:B12:H531	3:E:403:B12:H543	2.01	0.42
1:B:205:MET:HE1	3:B:403:B12:C43	2.49	0.42
3:B:403:B12:H552	3:B:403:B12:H531	2.01	0.42
1:E:130:ASP:OD1	3:E:403:B12:H542	2.20	0.42
1:B:234:TYR:CZ	1:B:284:GLY:HA2	2.54	0.42
1:D:92:ILE:CG2	1:D:260:ALA:O	2.67	0.42
1:D:191:LEU:HD21	1:D:201:GLY:HA2	2.01	0.42
1:A:191:LEU:HD21	1:A:201:GLY:HA2	2.02	0.42
1:B:294:ARG:HH11	1:B:294:ARG:CG	2.30	0.42
3:D:403:B12:H262	3:D:403:B12:H91	1.83	0.42
1:C:244:PRO:O	1:C:247:ARG:HB2	2.20	0.42
1:E:92:ILE:CG2	1:E:260:ALA:O	2.68	0.42
1:A:40:GLU:OE1	1:A:42:ARG:NH1	2.53	0.42
1:A:294:ARG:HH11	1:A:294:ARG:CG	2.27	0.42
1:E:195:PRO:HG3	1:E:238:ILE:HD13	2.02	0.42
1:E:206:ASN:C	1:E:206:ASN:ND2	2.77	0.42
1:C:54:ARG:HD2	1:C:133:ALA:O	2.20	0.42
1:B:200:ILE:HB	1:B:204:THR:HG22	2.01	0.41
1:C:99:LEU:HD23	1:C:103:TYR:HD2	1.84	0.41
1:E:40:GLU:OE1	1:E:42:ARG:NH1	2.50	0.41
1:A:55:ILE:O	1:A:57:PRO:HD3	2.20	0.41
1:A:92:ILE:HD11	1:A:96:SER:HB3	2.02	0.41
3:A:403:B12:H551	3:A:403:B12:C61	2.51	0.41
3:B:403:B12:C61	3:B:403:B12:H551	2.50	0.41
1:D:99:LEU:HD23	1:D:103:TYR:HD2	1.83	0.41
1:D:180:VAL:HG13	1:D:182:TYR:CD1	2.55	0.41
1:D:200:ILE:HB	1:D:204:THR:HG22	2.02	0.41
3:D:403:B12:H551	3:D:403:B12:O63	2.20	0.41
1:E:31:GLU:HG3	1:E:55:ILE:HD13	2.02	0.41
1:B:54:ARG:HD2	1:B:133:ALA:O	2.19	0.41
3:D:403:B12:H351	3:D:403:B12:C36	2.37	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:MET:HE1	3:A:403:B12:C43	2.51	0.41
3:B:403:B12:H91	3:B:403:B12:H262	1.72	0.41
1:C:180:VAL:HG13	1:C:182:TYR:CD1	2.55	0.41
3:D:403:B12:H253	3:D:403:B12:H301	1.93	0.41
1:E:30:LEU:O	1:E:31:GLU:C	2.64	0.41
1:E:54:ARG:HD2	1:E:133:ALA:O	2.19	0.41
1:B:206:ASN:C	1:B:206:ASN:ND2	2.75	0.41
1:C:55:ILE:O	1:C:57:PRO:HD3	2.21	0.41
1:E:244:PRO:O	1:E:247:ARG:HB2	2.20	0.41
1:B:135:VAL:HG12	1:B:137:THR:HG22	2.03	0.41
1:B:191:LEU:HD21	1:B:201:GLY:HA2	2.03	0.41
1:A:92:ILE:HG23	1:A:260:ALA:O	2.21	0.41
1:C:310:ALA:O	1:C:311:ILE:C	2.64	0.41
1:E:236:CYS:O	3:E:403:B12:O39	2.39	0.41
1:D:198:CYS:SG	2:D:402:SF4:S2	3.18	0.41
1:E:55:ILE:O	1:E:57:PRO:HD3	2.21	0.41
1:A:97:TRP:HB3	1:A:255:ILE:HD12	2.03	0.41
1:A:141:LYS:HZ3	3:A:403:B12:H4B	1.81	0.41
1:D:40:GLU:OE1	1:D:42:ARG:NH1	2.51	0.41
1:D:135:VAL:HG12	1:D:137:THR:HG22	2.03	0.41
1:D:244:PRO:O	1:D:247:ARG:HB2	2.21	0.41
1:D:54:ARG:HD2	1:D:133:ALA:O	2.21	0.41
1:E:135:VAL:HG12	1:E:137:THR:HG22	2.03	0.40
1:A:135:VAL:HG12	1:A:137:THR:HG22	2.03	0.40
1:D:55:ILE:O	1:D:57:PRO:HD3	2.21	0.40
1:D:92:ILE:HD13	1:D:92:ILE:HG21	1.75	0.40
1:E:99:LEU:HD23	1:E:103:TYR:CD2	2.56	0.40
1:A:195:PRO:HG3	1:A:238:ILE:HD13	2.04	0.40
1:C:31:GLU:HG3	1:C:55:ILE:HD13	2.03	0.40
1:C:137:THR:HG21	3:C:403:B12:H0R8	1.80	0.40
3:B:403:B12:H2R	3:B:403:B12:C7B	2.51	0.40
1:D:31:GLU:HG3	1:D:55:ILE:HD13	2.03	0.40
1:D:178:GLN:O	1:D:179:GLU:CB	2.69	0.40
1:D:195:PRO:HG3	1:D:238:ILE:HD13	2.04	0.40
3:D:403:B12:O7R	3:D:403:B12:C2B	2.69	0.40
1:B:31:GLU:HG3	1:B:55:ILE:HD13	2.03	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	366/399 (92%)	339 (93%)	24 (7%)	3 (1%)	16	27
1	B	366/399 (92%)	341 (93%)	21 (6%)	4 (1%)	11	19
1	C	363/399 (91%)	335 (92%)	25 (7%)	3 (1%)	16	27
1	D	366/399 (92%)	340 (93%)	22 (6%)	4 (1%)	11	19
1	E	366/399 (92%)	339 (93%)	24 (7%)	3 (1%)	16	27
All	All	1827/1995 (92%)	1694 (93%)	116 (6%)	17 (1%)	14	24

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62	GLU
1	A	179	GLU
1	B	62	GLU
1	B	179	GLU
1	B	252	PRO
1	C	62	GLU
1	C	179	GLU
1	C	252	PRO
1	D	62	GLU
1	D	179	GLU
1	D	252	PRO
1	E	62	GLU
1	E	179	GLU
1	E	252	PRO
1	A	252	PRO
1	D	208	ARG
1	B	208	ARG

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	314/342 (92%)	294 (94%)	20 (6%)	16	28
1	B	314/342 (92%)	294 (94%)	20 (6%)	16	28
1	C	311/342 (91%)	292 (94%)	19 (6%)	17	30
1	D	314/342 (92%)	296 (94%)	18 (6%)	18	32
1	E	314/342 (92%)	294 (94%)	20 (6%)	16	28
All	All	1567/1710 (92%)	1470 (94%)	97 (6%)	16	29

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ILE
1	A	18	LYS
1	A	28	ASP
1	A	31	GLU
1	A	54	ARG
1	A	55	ILE
1	A	65	LYS
1	A	80	GLN
1	A	92	ILE
1	A	99	LEU
1	A	129	VAL
1	A	137	THR
1	A	155	ILE
1	A	172	LEU
1	A	204	THR
1	A	208	ARG
1	A	250	ASP
1	A	294	ARG
1	A	310	VAL
1	A	334	GLU
1	B	3	ILE
1	B	18	LYS
1	B	28	ASP

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Mol	Chain	Res	Type
1	B	31	GLU
1	B	54	ARG
1	B	55	ILE
1	B	65	LYS
1	B	80	GLN
1	B	92	ILE
1	B	111	ARG
1	B	129	VAL
1	B	137	THR
1	B	155	ILE
1	B	204	THR
1	B	208	ARG
1	B	250	ASP
1	B	294	ARG
1	B	310	VAL
1	B	334	GLU
1	B	369	GLN
1	C	18	LYS
1	C	28	ASP
1	C	31	GLU
1	C	54	ARG
1	C	55	ILE
1	C	65	LYS
1	C	80	GLN
1	C	92	ILE
1	C	99	LEU
1	C	129	VAL
1	C	137	THR
1	C	155	ILE
1	C	172	LEU
1	C	204	THR
1	C	208	ARG
1	C	250	ASP
1	C	296	ARG
1	C	312	VAL
1	C	336	GLU
1	D	3	ILE
1	D	18	LYS
1	D	28	ASP
1	D	31	GLU
1	D	54	ARG
1	D	55	ILE

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Mol	Chain	Res	Type
1	D	65	LYS
1	D	80	GLN
1	D	92	ILE
1	D	129	VAL
1	D	137	THR
1	D	155	ILE
1	D	204	THR
1	D	208	ARG
1	D	250	ASP
1	D	294	ARG
1	D	310	VAL
1	D	334	GLU
1	E	3	ILE
1	E	18	LYS
1	E	28	ASP
1	E	31	GLU
1	E	54	ARG
1	E	55	ILE
1	E	65	LYS
1	E	80	GLN
1	E	92	ILE
1	E	99	LEU
1	E	129	VAL
1	E	137	THR
1	E	155	ILE
1	E	204	THR
1	E	208	ARG
1	E	250	ASP
1	E	294	ARG
1	E	310	VAL
1	E	334	GLU
1	E	369	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	HIS
1	A	50	ASN
1	A	80	GLN
1	A	95	ASN
1	A	178	GLN
1	A	206	ASN

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Mol	Chain	Res	Type
1	A	319	ASN
1	B	48	HIS
1	B	50	ASN
1	B	80	GLN
1	B	95	ASN
1	B	178	GLN
1	B	206	ASN
1	C	48	HIS
1	C	50	ASN
1	C	80	GLN
1	C	95	ASN
1	C	178	GLN
1	C	206	ASN
1	D	48	HIS
1	D	50	ASN
1	D	80	GLN
1	D	95	ASN
1	D	178	GLN
1	D	206	ASN
1	D	319	ASN
1	E	80	GLN
1	E	178	GLN
1	E	206	ASN
1	E	319	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

15 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$
2	SF4	B	401	1	0,12,12	-	-	-		
3	B12	C	403	-	94,101,101	0.77	3 (3%)	149,166,166	1.27	14 (9%)
3	B12	E	403	4	94,101,101	0.78	3 (3%)	149,166,166	1.25	16 (10%)
2	SF4	E	401	1	0,12,12	-	-	-		
3	B12	A	403	4	94,101,101	0.71	3 (3%)	149,166,166	1.37	20 (13%)
3	B12	B	403	-	94,101,101	1.04	4 (4%)	149,166,166	1.53	21 (14%)
2	SF4	B	402	1	0,12,12	-	-	-		
2	SF4	A	402	1	0,12,12	-	-	-		
2	SF4	E	402	1	0,12,12	-	-	-		
3	B12	D	403	-	94,101,101	0.84	4 (4%)	149,166,166	1.43	18 (12%)
2	SF4	D	402	1	0,12,12	-	-	-		
2	SF4	C	402	1	0,12,12	-	-	-		
2	SF4	D	401	1	0,12,12	-	-	-		
2	SF4	A	401	1	0,12,12	-	-	-		
2	SF4	C	401	1	0,12,12	-	-	-		

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
2	SF4	B	401	1	-	-	0/6/5/5
2	SF4	C	401	1	-	-	0/6/5/5
3	B12	E	403	4	-	11/56/223/223	0/3/11/11
2	SF4	E	401	1	-	-	0/6/5/5
3	B12	A	403	4	-	20/56/223/223	0/3/11/11
3	B12	B	403	-	-	17/56/223/223	0/3/11/11
2	SF4	B	402	1	-	-	0/6/5/5
2	SF4	A	402	1	-	-	0/6/5/5
2	SF4	E	402	1	-	-	0/6/5/5
3	B12	D	403	-	-	16/56/223/223	0/3/11/11
2	SF4	D	402	1	-	-	0/6/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SF4	C	402	1	-	-	0/6/5/5
2	SF4	D	401	1	-	-	0/6/5/5
2	SF4	A	401	1	-	-	0/6/5/5
3	B12	C	403	-	-	16/56/223/223	0/3/11/11

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	403	B12	C2B-N3B	-6.53	1.19	1.31
3	E	403	B12	C14-N23	4.78	1.41	1.35
3	C	403	B12	C2B-N3B	-4.78	1.22	1.31
3	D	403	B12	O3-C2P	-4.41	1.34	1.45
3	B	403	B12	O3-C2P	-4.00	1.35	1.45
3	D	403	B12	C2B-N3B	-3.90	1.24	1.31
3	B	403	B12	C11-N23	-3.81	1.29	1.36
3	B	403	B12	P-O3	-3.24	1.49	1.59
3	A	403	B12	O3-C2P	-3.20	1.37	1.45
3	E	403	B12	C11-N23	-3.16	1.31	1.36
3	A	403	B12	C11-N23	-2.74	1.31	1.36
3	C	403	B12	O3-C2P	-2.56	1.39	1.45
3	A	403	B12	P-O3	-2.31	1.52	1.59
3	E	403	B12	O3-C2P	-2.20	1.39	1.45
3	D	403	B12	C11-N23	-2.20	1.32	1.36
3	D	403	B12	P-O3	-2.19	1.52	1.59
3	C	403	B12	C14-N23	2.18	1.38	1.35

All (89) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	403	B12	C2P-C1P-N59	-7.67	101.64	112.92
3	D	403	B12	O2-P-O3	-7.42	82.16	102.87
3	B	403	B12	C9-N22-C6	6.43	113.02	105.28
3	B	403	B12	O2-P-O3	-6.15	85.70	102.87
3	B	403	B12	C8B-N1B-C1R	-5.99	115.28	125.41
3	C	403	B12	O2-P-O3	-5.57	87.30	102.87
3	A	403	B12	C8B-N1B-C1R	-5.49	116.13	125.41
3	E	403	B12	C3P-C2P-C1P	-4.68	102.38	111.42
3	E	403	B12	C2P-C1P-N59	-4.56	106.22	112.92
3	C	403	B12	C41-C8-C9	-4.52	103.31	111.19
3	A	403	B12	C9-N22-C6	4.44	110.62	105.28
3	E	403	B12	C9-N22-C6	4.23	110.37	105.28
3	B	403	B12	C36-C7-C37	4.07	117.66	110.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	403	B12	O2-P-O3	-4.03	91.61	102.87
3	B	403	B12	C2P-C1P-N59	-3.70	107.48	112.92
3	A	403	B12	O2-P-O4	3.67	121.53	109.81
3	C	403	B12	C19-N24-C16	3.43	111.03	107.29
3	D	403	B12	C9B-N3B-C2B	3.35	108.51	104.40
3	D	403	B12	C9-N22-C6	3.28	109.23	105.28
3	C	403	B12	C9-N22-C6	3.20	109.13	105.28
3	B	403	B12	C10-C9-N22	3.18	129.37	125.74
3	B	403	B12	C30-C3-C2	-3.17	112.02	119.00
3	E	403	B12	C55-C56-C57	-3.15	104.22	111.25
3	B	403	B12	C8B-N1B-C2B	-3.12	103.43	106.26
3	B	403	B12	N1B-C2B-N3B	3.10	118.33	113.94
3	B	403	B12	C13-C14-C15	-3.08	119.63	124.32
3	E	403	B12	C41-C8-C9	-3.06	105.85	111.19
3	A	403	B12	C56-C55-C17	-3.03	109.74	115.58
3	E	403	B12	O2-P-O3	-2.99	94.51	102.87
3	A	403	B12	C30-C3-C2	-2.99	112.41	119.00
3	C	403	B12	C36-C7-C37	2.89	115.64	110.74
3	A	403	B12	C15-C16-N24	-2.88	118.32	122.42
3	A	403	B12	C15-C14-N23	2.87	129.73	126.26
3	D	403	B12	C55-C56-C57	-2.86	104.87	111.25
3	A	403	B12	C19-C1-N21	-2.83	99.23	102.14
3	D	403	B12	C10-C9-N22	2.79	128.93	125.74
3	A	403	B12	C2P-C1P-N59	-2.78	108.84	112.92
3	A	403	B12	O3-C2P-C1P	-2.78	101.44	106.94
3	D	403	B12	C56-C55-C17	-2.76	110.27	115.58
3	C	403	B12	O3-P-O4	2.76	118.61	109.81
3	C	403	B12	O2-P-O4	2.75	118.58	109.81
3	C	403	B12	C55-C56-C57	-2.74	105.14	111.25
3	A	403	B12	C26-C2-C3	-2.73	102.66	107.42
3	B	403	B12	C15-C14-N23	2.72	129.54	126.26
3	D	403	B12	C54-C17-C55	-2.70	104.77	109.27
3	E	403	B12	C8B-N1B-C1R	-2.69	120.85	125.41
3	D	403	B12	O2-P-O4	2.68	118.38	109.81
3	B	403	B12	C15-C16-N24	-2.64	118.66	122.42
3	D	403	B12	C8B-N1B-C1R	-2.64	120.94	125.41
3	E	403	B12	O3-P-O4	2.63	118.21	109.81
3	C	403	B12	C1-C19-N24	-2.62	103.34	106.25
3	A	403	B12	C10-C9-N22	2.60	128.71	125.74
3	B	403	B12	C19-N24-C16	2.60	110.13	107.29
3	D	403	B12	C26-C2-C3	-2.60	102.89	107.42
3	C	403	B12	C9B-N3B-C2B	2.59	107.58	104.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	403	B12	C18-C19-N24	2.59	106.22	102.33
3	E	403	B12	C15-C16-N24	-2.59	118.74	122.42
3	A	403	B12	C55-C17-C18	-2.57	106.21	111.12
3	A	403	B12	C20-C1-C2	2.53	117.46	113.28
3	B	403	B12	C55-C56-C57	-2.48	105.72	111.25
3	B	403	B12	O2-C3R-C2R	-2.46	102.85	111.68
3	A	403	B12	C20-C1-C19	-2.46	106.98	109.35
3	B	403	B12	C55-C17-C18	-2.45	106.45	111.12
3	E	403	B12	C17-C16-N24	2.44	114.89	111.17
3	C	403	B12	C1-C19-C18	-2.42	117.98	121.90
3	D	403	B12	O5-P-O3	2.41	116.53	106.70
3	A	403	B12	C26-C2-C1	2.41	113.73	110.00
3	E	403	B12	C30-C3-C2	-2.40	113.71	119.00
3	D	403	B12	C13-C14-C15	-2.37	120.71	124.32
3	E	403	B12	C15-C14-N23	2.37	129.12	126.26
3	B	403	B12	C1-C19-C18	-2.35	118.09	121.90
3	A	403	B12	C36-C7-C37	2.30	114.65	110.74
3	E	403	B12	C20-C1-C2	2.29	117.07	113.28
3	C	403	B12	C2P-C1P-N59	-2.27	109.58	112.92
3	D	403	B12	C19-N24-C16	2.24	109.74	107.29
3	D	403	B12	C17-C16-N24	2.22	114.55	111.17
3	A	403	B12	C13-C14-C15	-2.22	120.94	124.32
3	C	403	B12	C10-C9-N22	2.21	128.26	125.74
3	B	403	B12	O3-C2P-C1P	-2.20	102.58	106.94
3	A	403	B12	C19-N24-C16	2.14	109.63	107.29
3	E	403	B12	C26-C2-C3	-2.13	103.70	107.42
3	B	403	B12	O5-P-O3	2.11	115.27	106.70
3	C	403	B12	C19-C1-N21	-2.10	99.97	102.14
3	E	403	B12	O5-P-O2	2.08	115.18	106.70
3	B	403	B12	O7R-C2R-C3R	-2.08	105.35	111.19
3	B	403	B12	C54-C17-C55	-2.06	105.85	109.27
3	D	403	B12	C53-C15-C16	-2.03	116.92	120.36
3	D	403	B12	C8B-C9B-N3B	-2.03	107.80	110.00
3	D	403	B12	C9-C10-C11	-2.03	123.07	125.97

There are no chirality outliers.

All (80) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	403	B12	C4-C3-C30-C31
3	A	403	B12	C56-C57-N59-C1P
3	A	403	B12	N59-C1P-C2P-C3P

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Mol	Chain	Res	Type	Atoms
3	A	403	B12	N59-C1P-C2P-O3
3	A	403	B12	C3R-O2-P-O3
3	A	403	B12	C3R-O2-P-O5
3	B	403	B12	N59-C1P-C2P-C3P
3	B	403	B12	N59-C1P-C2P-O3
3	B	403	B12	C1P-C2P-O3-P
3	B	403	B12	C3P-C2P-O3-P
3	B	403	B12	C3R-O2-P-O3
3	B	403	B12	C3R-O2-P-O5
3	B	403	B12	C2R-C1R-N1B-C2B
3	C	403	B12	N59-C1P-C2P-C3P
3	C	403	B12	N59-C1P-C2P-O3
3	C	403	B12	C2P-O3-P-O4
3	C	403	B12	C2P-O3-P-O2
3	C	403	B12	C3R-O2-P-O3
3	D	403	B12	C56-C57-N59-C1P
3	D	403	B12	N59-C1P-C2P-O3
3	D	403	B12	C2P-O3-P-O2
3	D	403	B12	C3R-O2-P-O3
3	E	403	B12	C56-C57-N59-C1P
3	E	403	B12	N59-C1P-C2P-C3P
3	E	403	B12	N59-C1P-C2P-O3
3	E	403	B12	C3P-C2P-O3-P
3	E	403	B12	C3R-O2-P-O3
3	A	403	B12	O58-C57-N59-C1P
3	D	403	B12	O58-C57-N59-C1P
3	E	403	B12	O58-C57-N59-C1P
3	E	403	B12	C3R-C4R-C5R-O8R
3	E	403	B12	O6R-C4R-C5R-O8R
3	A	403	B12	C2-C3-C30-C31
3	A	403	B12	C42-C41-C8-C9
3	B	403	B12	C56-C57-N59-C1P
3	A	403	B12	C3R-C4R-C5R-O8R
3	B	403	B12	C3-C2-C26-C27
3	E	403	B12	C17-C18-C60-C61
3	A	403	B12	O6R-C4R-C5R-O8R
3	B	403	B12	O58-C57-N59-C1P
3	D	403	B12	C3R-C4R-C5R-O8R
3	D	403	B12	C4-C3-C30-C31
3	D	403	B12	O6R-C4R-C5R-O8R
3	A	403	B12	C42-C41-C8-C7
3	B	403	B12	C2-C3-C30-C31

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Mol	Chain	Res	Type	Atoms
3	C	403	B12	C12-C13-C48-C49
3	D	403	B12	C2-C3-C30-C31
3	B	403	B12	C2R-C1R-N1B-C8B
3	C	403	B12	C2R-C1R-N1B-C8B
3	C	403	B12	O6R-C1R-N1B-C8B
3	C	403	B12	C56-C57-N59-C1P
3	C	403	B12	O58-C57-N59-C1P
3	A	403	B12	C8-C41-C42-C43
3	B	403	B12	C4-C3-C30-C31
3	C	403	B12	C1P-C2P-O3-P
3	B	403	B12	C25-C2-C26-C27
3	B	403	B12	C7-C37-C38-N40
3	A	403	B12	C2P-O3-P-O4
3	B	403	B12	C3R-O2-P-O4
3	C	403	B12	C3R-O2-P-O4
3	D	403	B12	C2P-O3-P-O5
3	E	403	B12	C3R-O2-P-O5
3	A	403	B12	C12-C13-C48-C49
3	A	403	B12	C2P-O3-P-O2
3	D	403	B12	N59-C1P-C2P-C3P
3	A	403	B12	C3-C30-C31-C32
3	B	403	B12	C7-C37-C38-O39
3	A	403	B12	C2P-C1P-N59-C57
3	C	403	B12	C3P-C2P-O3-P
3	D	403	B12	C1P-C2P-O3-P
3	C	403	B12	C2P-C1P-N59-C57
3	D	403	B12	C8-C41-C42-C43
3	A	403	B12	C3R-O2-P-O4
3	C	403	B12	C3R-O2-P-O5
3	D	403	B12	C2P-O3-P-O4
3	D	403	B12	C3R-O2-P-O4
3	D	403	B12	C3R-O2-P-O5
3	E	403	B12	C3R-O2-P-O4
3	A	403	B12	C1P-C2P-O3-P
3	C	403	B12	C16-C17-C55-C56

There are no ring outliers.

9 monomers are involved in 105 short contacts:

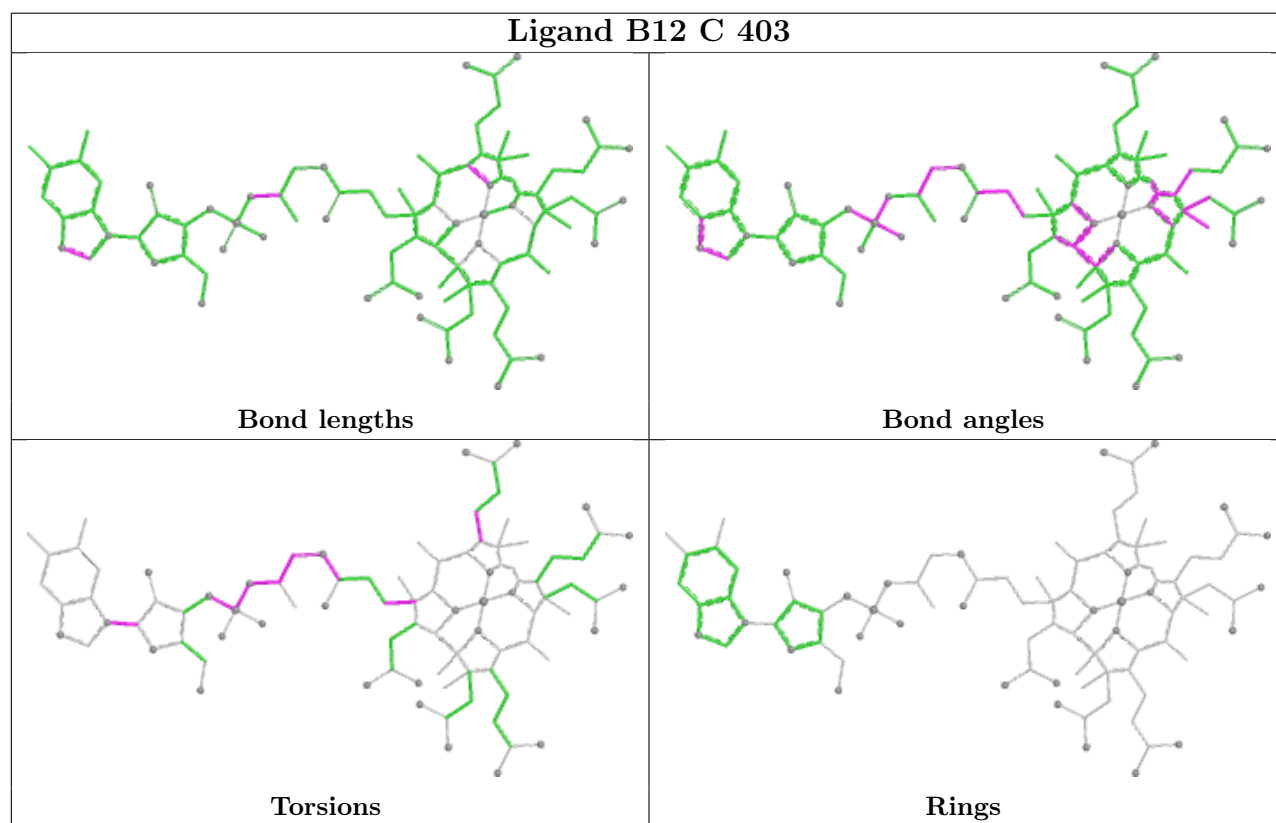
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	403	B12	16	0
3	E	403	B12	22	0

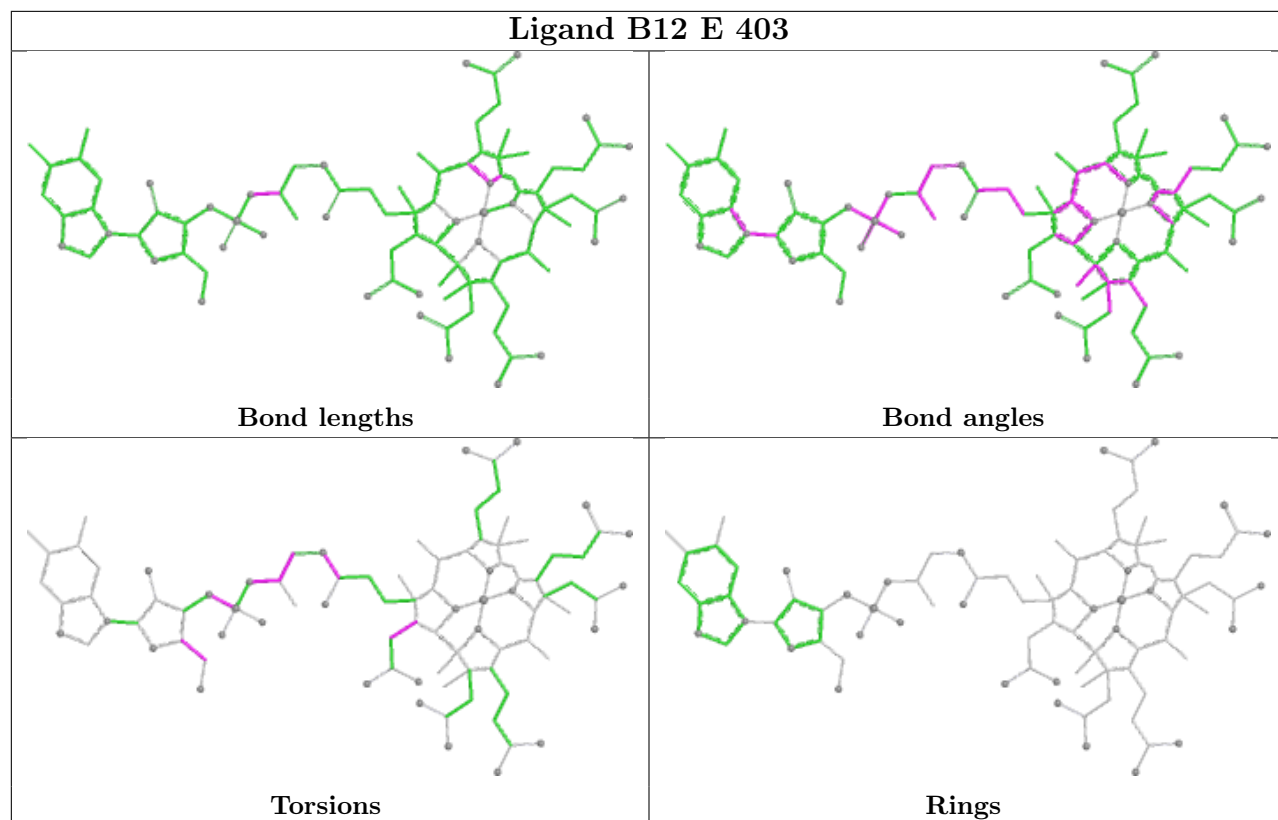
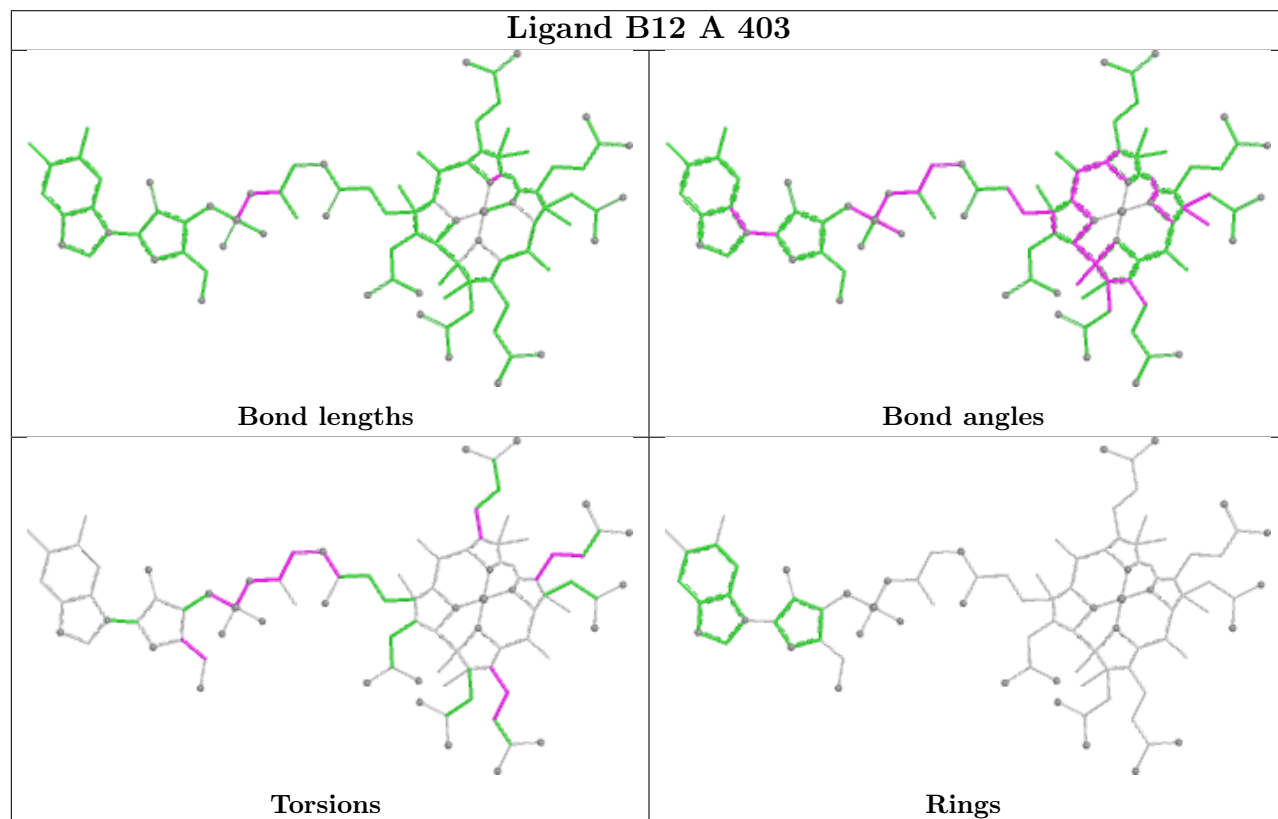
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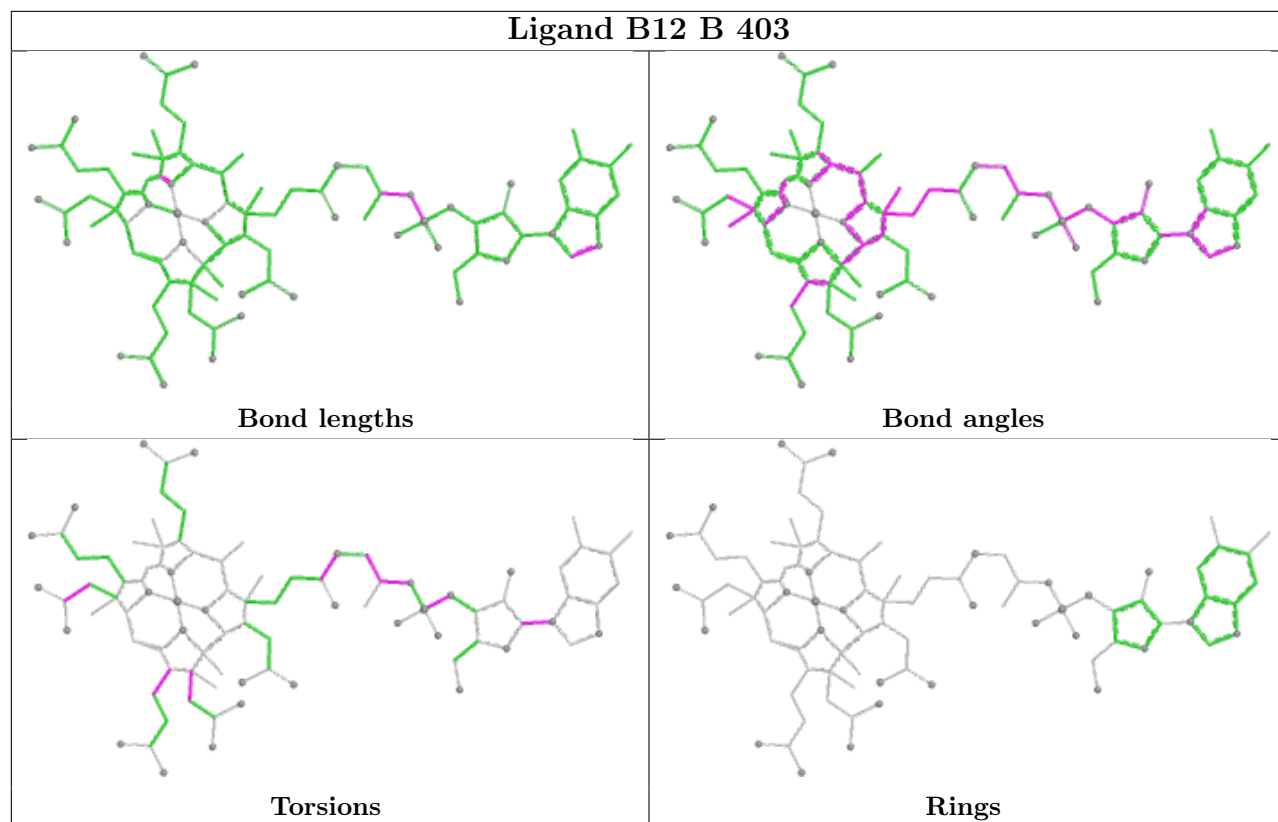
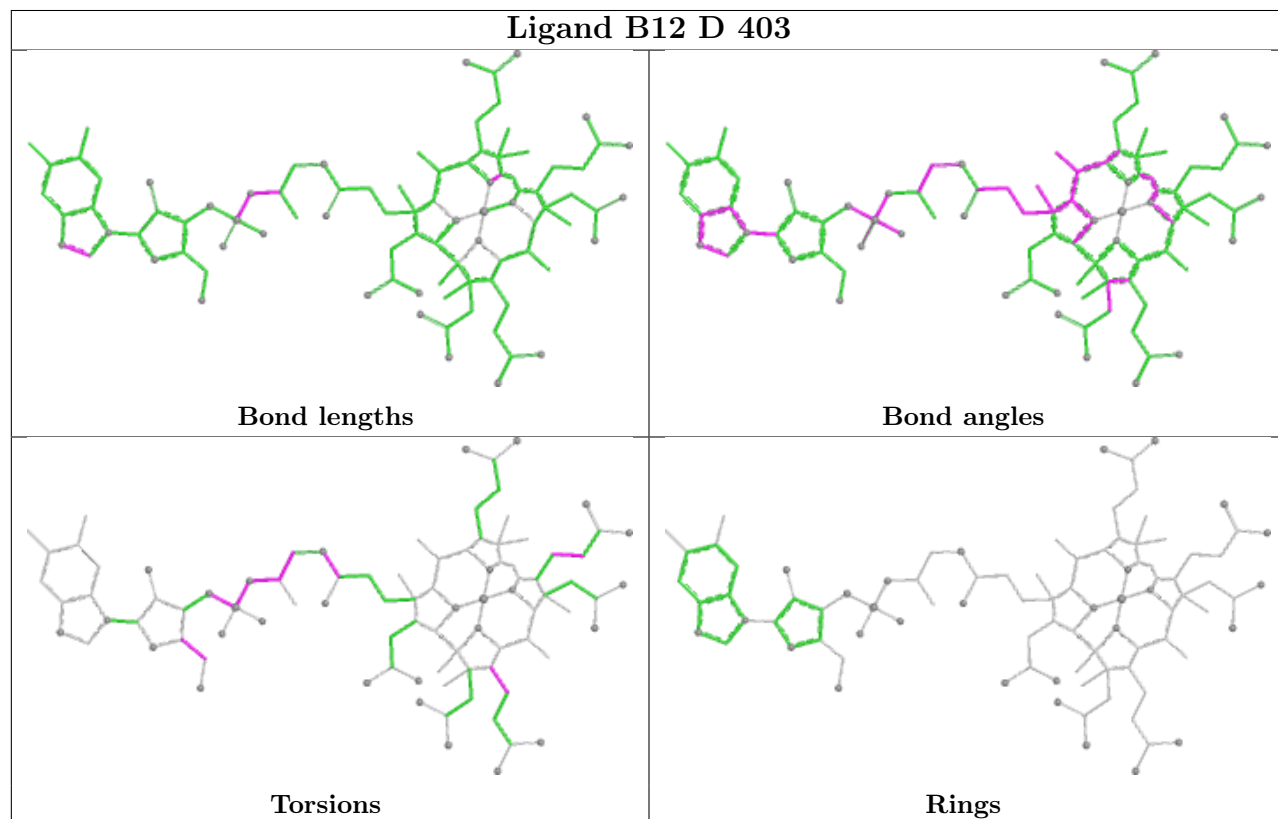
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	403	B12	24	0
3	B	403	B12	22	0
2	B	402	SF4	1	0
2	A	402	SF4	1	0
2	E	402	SF4	1	0
3	D	403	B12	16	0
2	D	402	SF4	2	0

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 B12 E 403**Ligand B12 A 403**

Ligand B12 B 403**Ligand B12 D 403**

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	370/399 (92%)	-0.04	3 (0%) 82 79	36, 67, 103, 130	0
1	B	370/399 (92%)	-0.03	3 (0%) 82 79	39, 66, 108, 169	0
1	C	367/399 (91%)	0.33	14 (3%) 44 36	41, 78, 122, 148	0
1	D	370/399 (92%)	0.14	4 (1%) 78 75	46, 75, 116, 158	0
1	E	370/399 (92%)	0.33	8 (2%) 62 56	45, 82, 131, 181	0
All	All	1847/1995 (92%)	0.15	32 (1%) 69 64	36, 73, 119, 181	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	253	LEU	4.4
1	D	253	LEU	4.3
1	C	253	LEU	3.3
1	E	112	LEU	3.2
1	B	103	TYR	3.2
1	C	145	ILE	3.1
1	B	253	LEU	3.0
1	A	1	MET	2.9
1	C	163	MET	2.9
1	E	69	SER	2.8
1	A	103	TYR	2.7
1	C	355	LYS	2.7
1	A	253	LEU	2.6
1	C	164	TYR	2.6
1	D	248	GLY	2.6
1	B	353	LYS	2.5
1	C	249	ILE	2.5
1	E	109	LEU	2.5
1	C	184	CYS	2.3
1	E	248	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	11	ALA	2.2
1	C	71	ALA	2.2
1	C	181	ASP	2.2
1	C	245	TYR	2.1
1	C	20	GLY	2.1
1	C	137	THR	2.1
1	C	5	LEU	2.1
1	C	168	LEU	2.1
1	D	119	LEU	2.1
1	E	71	ALA	2.0
1	E	3	ILE	2.0
1	D	103	TYR	2.0

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(Å ²)	Q<0.9
3	B12	C	403	91/91	0.95	0.10	52,71,102,110	0
3	B12	E	403	91/91	0.95	0.09	54,74,103,124	0
2	SF4	D	401	8/8	0.96	0.06	70,85,91,92	0
3	B12	A	403	91/91	0.96	0.10	42,55,108,118	0
3	B12	B	403	91/91	0.96	0.10	37,52,118,135	0
2	SF4	C	401	8/8	0.96	0.06	82,101,105,107	0
3	B12	D	403	91/91	0.96	0.10	41,59,106,122	0
2	SF4	C	402	8/8	0.96	0.06	62,72,75,76	0
2	SF4	E	402	8/8	0.97	0.05	69,74,77,82	0
2	SF4	E	401	8/8	0.97	0.06	89,102,106,110	0
2	SF4	A	401	8/8	0.98	0.04	65,73,76,78	0

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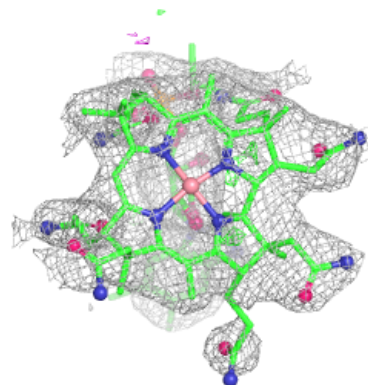
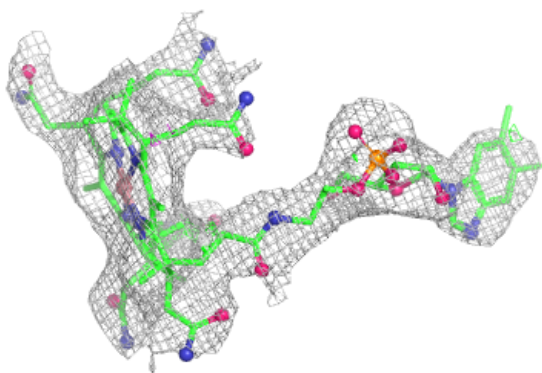
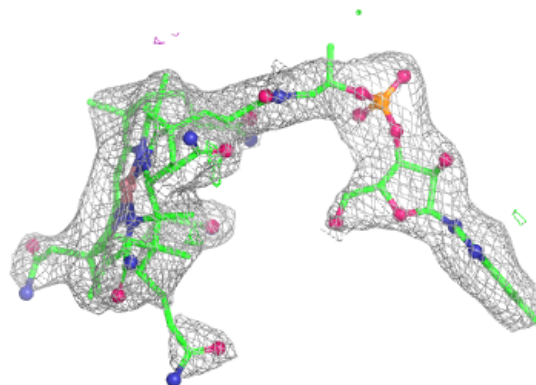
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SF4	B	401	8/8	0.98	0.04	63,69,75,76	0
2	SF4	A	402	8/8	0.99	0.04	53,57,60,64	0
2	SF4	D	402	8/8	0.99	0.04	51,58,62,66	0
2	SF4	B	402	8/8	0.99	0.03	46,51,55,55	0

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.

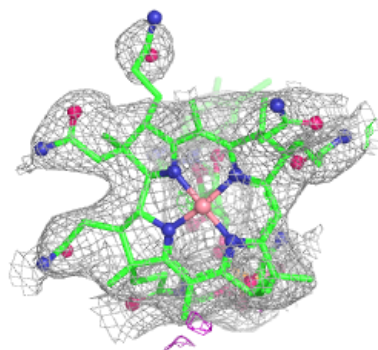
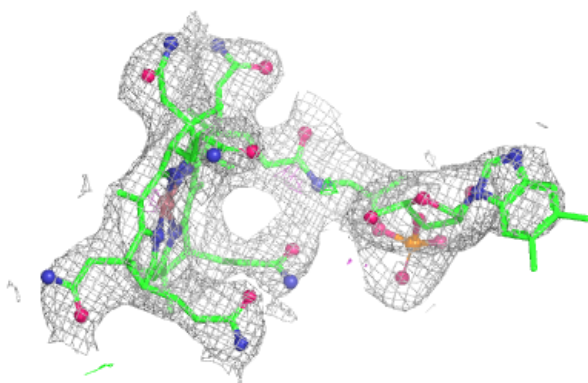
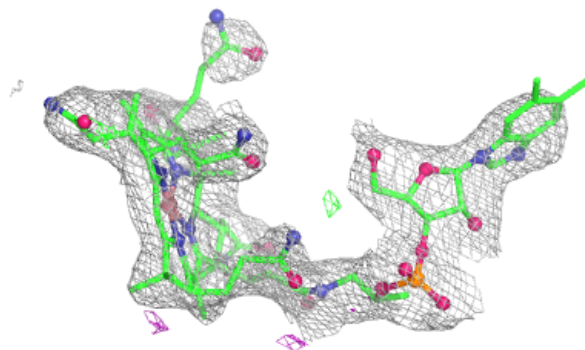
Electron density around B12 C 403:

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

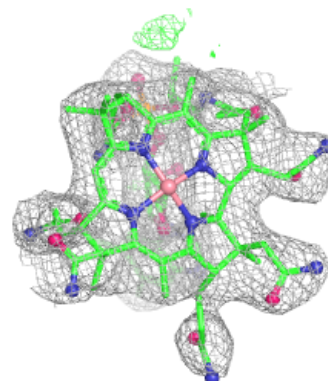
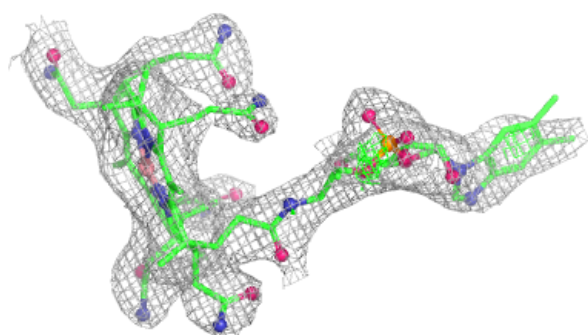
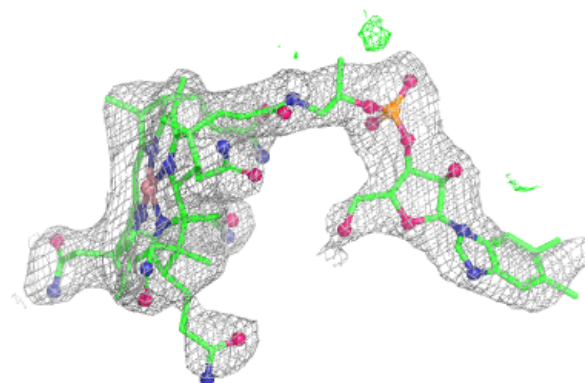


Electron density around B12 E 403:

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

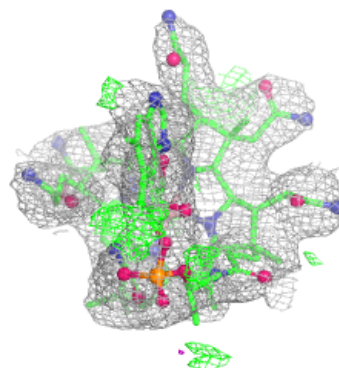
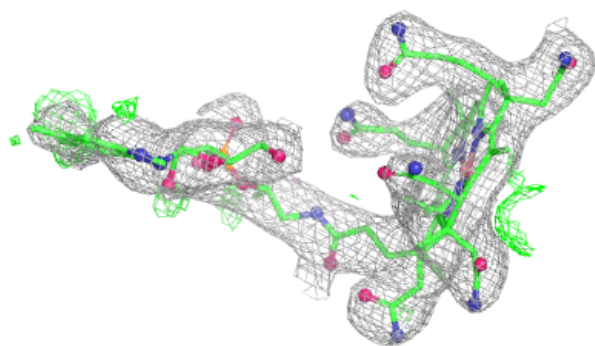
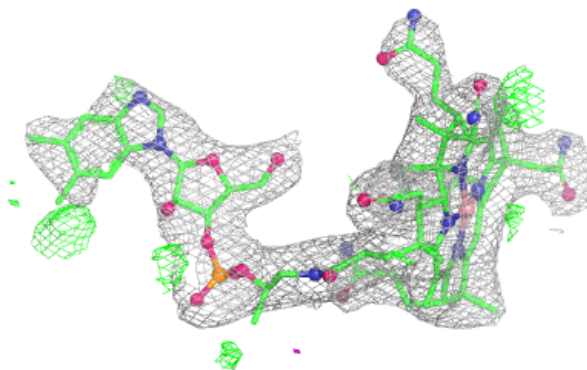
**Electron density around B12 A 403:**

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

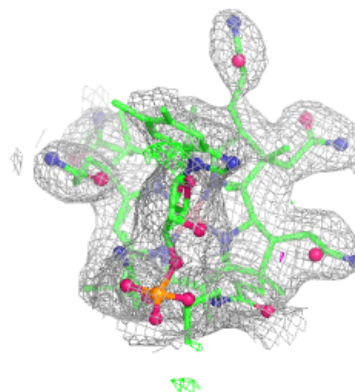
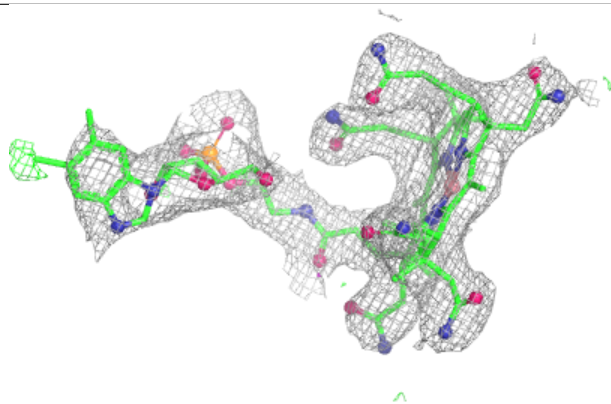
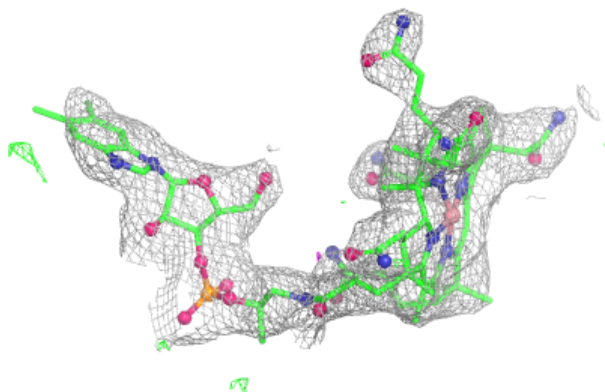


Electron density around B12 B 403:

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

**Electron density around B12 D 403:**

$2mF_o-DF_c$ (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.