



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:26 PM UTC

PDB ID : 3IEL / pdb_00003iel
Title : Crystal Structure of TTHA0252 from *Thermus thermophilus* HB8 complexed with UMP
Authors : Ishikawa, H.; Nakagawa, N.; Kuramitsu, S.; Yokoyama, S.; Masui, R.
Deposited on : 2009-07-23
Resolution : 2.35 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

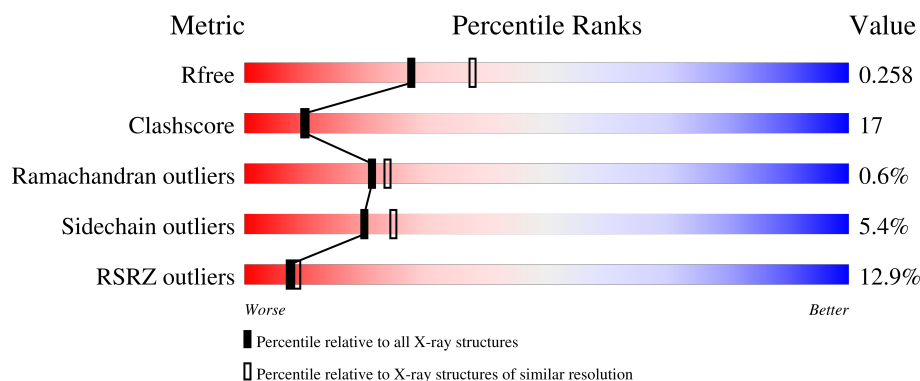
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	431	2% 68% 28% .
1	B	431	3% 71% 27% ..
1	C	431	19% 58% 40% .
1	D	431	29% 56% 41% .

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	SO4	A	433	-	-	X	-
2	SO4	A	443	-	-	X	-
2	SO4	C	453	-	-	X	-

2 Entry composition [i](#)

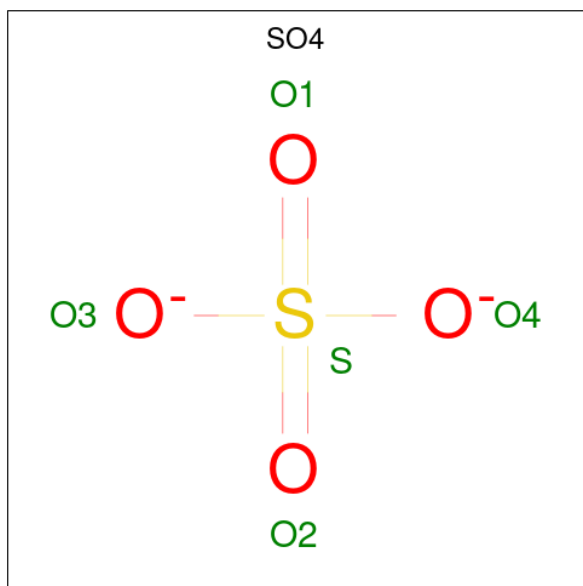
There are 6 unique types of molecules in this entry. The entry contains 14251 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Ribonuclease TTHA0252.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	431	Total	C	N	O	S	0	0	0
			3326	2127	597	594	8			
1	B	431	Total	C	N	O	S	0	0	0
			3326	2127	597	594	8			
1	C	431	Total	C	N	O	S	0	0	0
			3326	2127	597	594	8			
1	D	431	Total	C	N	O	S	0	0	0
			3326	2127	597	594	8			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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

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

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

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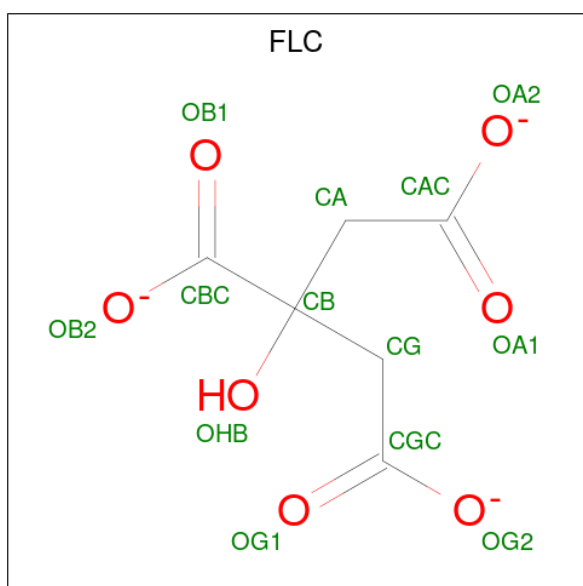
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total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

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

- Molecule 3 is CITRATE ANION (CCD ID: FLC) (formula: $C_6H_5O_7^-$).



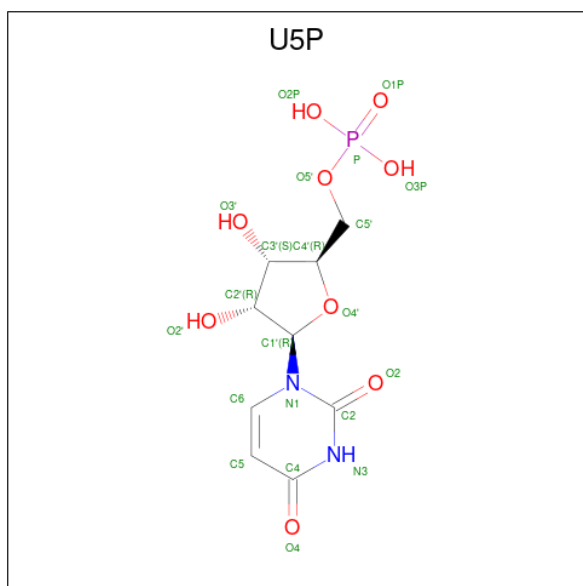
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	6	7		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			13	6	7		

- Molecule 4 is URIDINE-5'-MONOPHOSPHATE (CCD ID: U5P) (formula: $C_9H_{13}N_2O_9P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			21	9	2	9	1		
4	A	1	Total	C	N	O	P	0	0
			21	9	2	9	1		
4	A	1	Total	C	N	O	P	0	0
			21	9	2	9	1		
4	B	1	Total	C	N	O	P	0	0
			21	9	2	9	1		
4	B	1	Total	C	N	O	P	0	0
			21	9	2	9	1		
4	C	1	Total	C	N	O	P	0	0
			21	9	2	9	1		
4	D	1	Total	C	N	O	P	0	0
			21	9	2	9	1		

- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Zn	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total 2	Zn 2	0	0
5	C	2	Total 2	Zn 2	0	0
5	D	2	Total 2	Zn 2	0	0

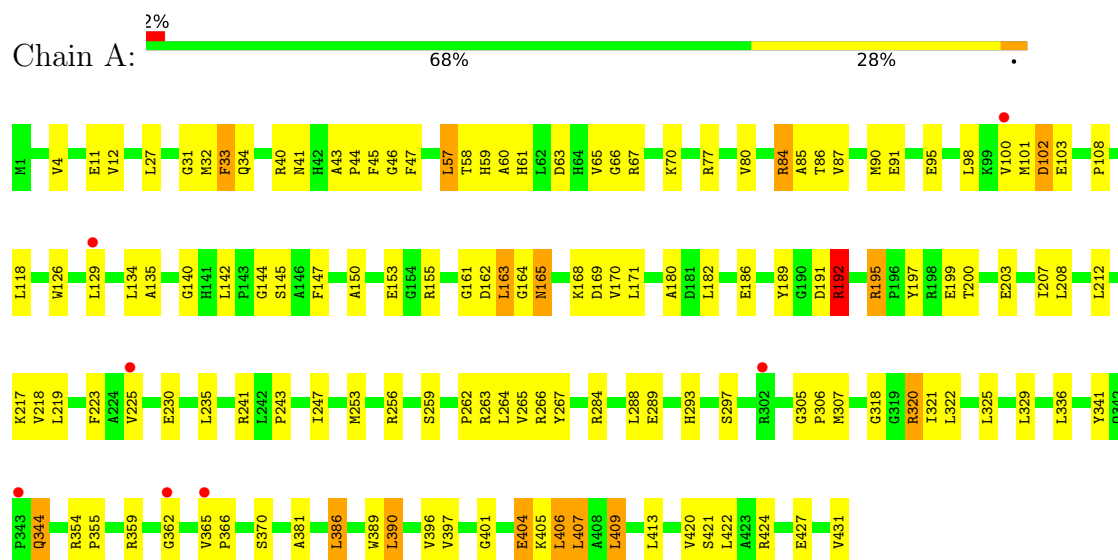
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	118	Total 118	O 118	0	0
6	B	95	Total 95	O 95	0	0
6	C	47	Total 47	O 47	0	0
6	D	26	Total 26	O 26	0	0

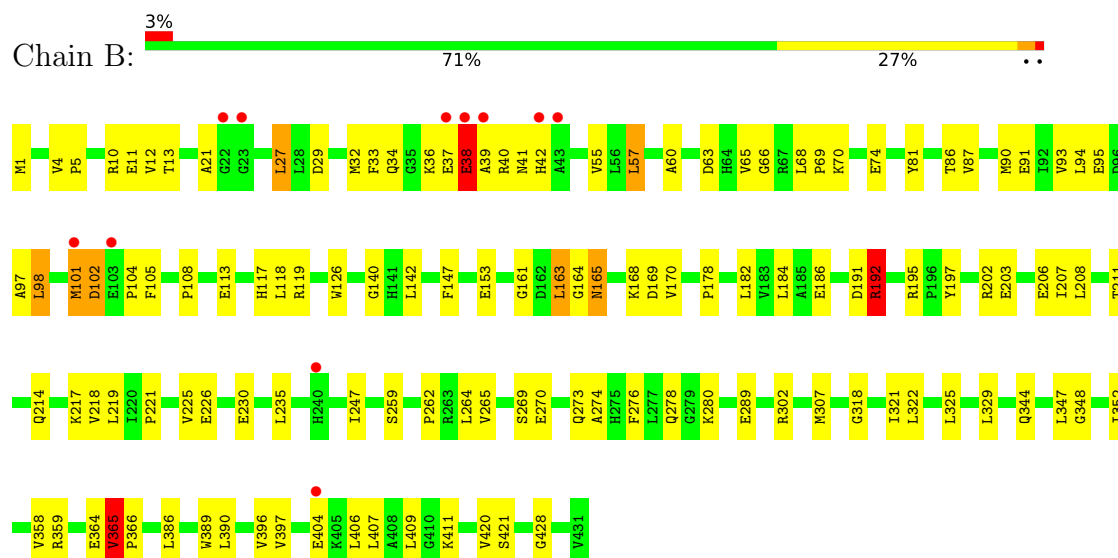
3 Residue-property plots

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: Ribonuclease TTHA0252

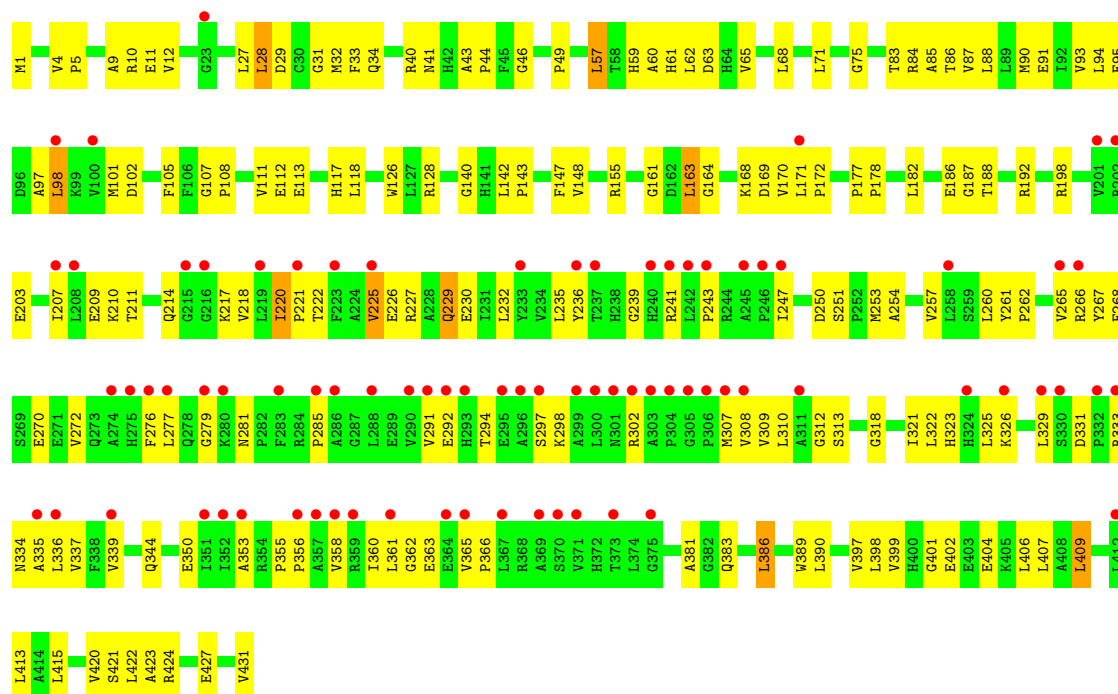


• Molecule 1: Ribonuclease TTHA0252

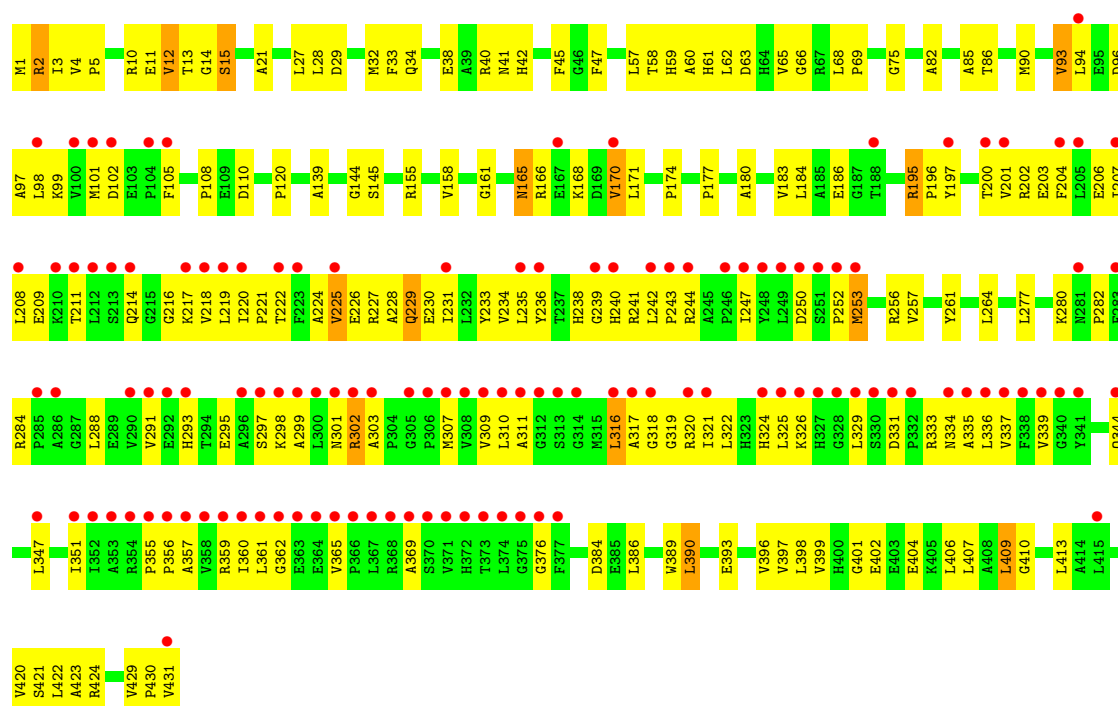


• Molecule 1: Ribonuclease TTHA0252





• Molecule 1: Ribonuclease TTHA0252



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	145.24Å 146.68Å 120.35Å 90.00° 110.11° 90.00°	Depositor
Resolution (Å)	50.00 – 2.35 50.00 – 2.35	Depositor EDS
% Data completeness (in resolution range)	95.3 (50.00-2.35) 95.3 (50.00-2.35)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.28 (at 2.34Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.222 , 0.259 0.222 , 0.258	Depositor DCC
R_{free} test set	9797 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	45.9	Xtriage
Anisotropy	0.257	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 49.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14251	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% 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: U5P, FLC, ZN, SO4

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	0.45	0/3407	0.99	19/4621 (0.4%)
1	B	0.45	0/3407	0.98	14/4621 (0.3%)
1	C	0.38	0/3407	0.92	13/4621 (0.3%)
1	D	0.34	0/3407	0.89	11/4621 (0.2%)
All	All	0.41	0/13628	0.95	57/18484 (0.3%)

There are no bond length outliers.

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	192	ARG	N-CA-C	11.99	118.92	108.13
1	A	192	ARG	N-CA-C	11.90	118.93	108.22
1	A	225	VAL	N-CA-C	11.47	121.31	110.53
1	B	225	VAL	N-CA-C	9.82	120.64	110.72
1	A	170	VAL	N-CA-C	8.00	118.80	110.72
1	D	75	GLY	N-CA-C	7.26	122.28	113.79
1	C	265	VAL	N-CA-C	7.11	117.82	110.36
1	A	100	VAL	N-CA-C	6.94	113.64	106.21
1	A	169	ASP	N-CA-C	6.82	121.73	113.41
1	B	195	ARG	N-CA-C	-6.79	100.72	110.08
1	B	60	ALA	N-CA-C	6.75	121.27	112.89
1	C	75	GLY	N-CA-C	6.70	120.75	113.58
1	B	170	VAL	N-CA-C	6.62	119.36	111.09
1	C	161	GLY	N-CA-C	-6.58	100.06	111.50
1	A	147	PHE	N-CA-C	-6.56	99.97	110.14
1	C	60	ALA	N-CA-C	6.49	120.94	112.89
1	B	147	PHE	N-CA-C	-6.22	100.03	109.85
1	B	66	GLY	N-CA-C	6.20	120.39	112.83
1	D	161	GLY	N-CA-C	-6.18	100.53	112.02
1	B	169	ASP	N-CA-C	6.16	120.93	113.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	4	VAL	N-CA-C	6.15	114.18	107.60
1	D	4	VAL	N-CA-C	6.07	113.56	107.55
1	C	148	VAL	N-CA-C	6.06	116.60	108.11
1	C	169	ASP	N-CA-C	6.05	122.12	114.31
1	D	12	VAL	N-CA-C	6.03	117.97	111.58
1	A	161	GLY	N-CA-C	-5.99	100.90	111.80
1	A	60	ALA	N-CA-C	5.97	120.18	113.01
1	B	161	GLY	N-CA-C	-5.92	101.03	111.80
1	C	46	GLY	N-CA-C	-5.90	106.38	114.64
1	C	225	VAL	N-CA-C	5.81	116.02	110.74
1	A	4	VAL	N-CA-C	5.71	114.24	107.73
1	B	365	VAL	N-CA-C	5.60	116.56	107.71
1	D	66	GLY	N-CA-C	5.59	119.65	112.83
1	D	60	ALA	N-CA-C	5.58	119.35	112.54
1	A	33	PHE	N-CA-C	-5.55	101.50	110.32
1	B	365	VAL	CA-C-N	5.52	125.79	119.83
1	B	365	VAL	C-N-CA	5.52	125.79	119.83
1	A	293	HIS	N-CA-C	5.45	118.46	109.85
1	D	93	VAL	N-CA-C	5.43	115.53	110.42
1	A	195	ARG	N-CA-C	-5.43	102.75	109.64
1	D	170	VAL	N-CA-C	5.40	115.60	110.53
1	A	66	GLY	N-CA-C	5.39	120.30	113.24
1	A	305	GLY	N-CA-C	-5.37	105.73	112.23
1	A	70	LYS	N-CA-C	-5.34	105.63	111.82
1	A	12	VAL	N-CA-C	5.25	120.26	113.07
1	A	223	PHE	N-CA-C	-5.25	102.72	110.48
1	D	317	ALA	N-CA-C	-5.09	106.90	113.01
1	B	184	LEU	N-CA-C	-5.09	100.22	108.41
1	A	180	ALA	N-CA-C	5.08	117.10	109.23
1	B	93	VAL	N-CA-C	5.07	115.29	110.42
1	C	220	ILE	N-CA-C	5.06	113.16	108.15
1	A	100	VAL	CB-CA-C	-5.06	107.55	113.22
1	C	107	GLY	CA-C-N	5.05	124.66	119.05
1	C	107	GLY	C-N-CA	5.05	124.66	119.05
1	D	177	PRO	CA-C-N	-5.05	115.52	120.52
1	D	177	PRO	C-N-CA	-5.05	115.52	120.52
1	C	147	PHE	N-CA-C	-5.03	102.75	110.14

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	3326	0	3351	96	0
1	B	3326	0	3351	84	0
1	C	3326	0	3351	144	0
1	D	3326	0	3351	159	0
2	A	160	0	0	9	0
2	B	115	0	0	0	0
2	C	125	0	0	5	0
2	D	80	0	0	1	0
3	A	13	0	5	1	0
3	B	13	0	5	2	0
4	A	63	0	33	2	0
4	B	42	0	22	2	0
4	C	21	0	11	0	0
4	D	21	0	11	0	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
5	C	2	0	0	0	0
5	D	2	0	0	0	0
6	A	118	0	0	5	0
6	B	95	0	0	6	0
6	C	47	0	0	3	0
6	D	26	0	0	1	0
All	All	14251	0	13491	476	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 17.

All (476) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:98:LEU:HD11	1:C:108:PRO:HA	1.40	1.03
1:D:235:LEU:HD23	1:D:247:ILE:HD13	1.37	1.03
1:B:33:PHE:H	1:B:41:ASN:HD21	1.05	1.01
1:D:33:PHE:H	1:D:41:ASN:HD21	1.10	0.99
1:B:153:GLU:HG2	6:B:519:HOH:O	1.65	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:GLU:HG2	2:A:433:SO4:O1	1.63	0.97
1:A:33:PHE:H	1:A:41:ASN:HD21	1.05	0.94
1:A:306:PRO:HD2	1:B:302:ARG:HH22	1.33	0.91
1:C:33:PHE:H	1:C:41:ASN:HD21	1.15	0.89
1:B:265:VAL:HG12	6:B:532:HOH:O	1.76	0.85
1:D:222:THR:HG22	1:D:339:VAL:HG21	1.58	0.83
1:D:284:ARG:HD3	1:D:288:LEU:HD23	1.59	0.83
1:D:1:MET:HG3	1:D:21:ALA:HB2	1.59	0.82
1:D:86:THR:HG22	1:D:90:MET:HE2	1.61	0.81
1:A:208:LEU:HD23	1:A:218:VAL:HG21	1.63	0.81
1:C:360:ILE:HG22	1:C:361:LEU:HD13	1.63	0.80
1:A:153:GLU:HA	2:A:433:SO4:O3	1.82	0.80
1:A:404:GLU:CD	1:A:404:GLU:H	1.89	0.80
1:D:98:LEU:HD21	1:D:108:PRO:HB3	1.64	0.80
1:A:168:LYS:HE3	1:A:230:GLU:OE1	1.81	0.79
1:D:168:LYS:HE2	1:D:230:GLU:OE1	1.83	0.79
1:B:86:THR:HG22	1:B:90:MET:HE2	1.64	0.79
1:B:97:ALA:O	1:B:101:MET:HB2	1.81	0.78
1:D:322:LEU:HB3	1:D:361:LEU:HD11	1.64	0.78
1:D:211:THR:HG21	1:D:335:ALA:HB2	1.64	0.78
1:B:101:MET:HG2	1:B:104:PRO:HB3	1.68	0.75
1:C:168:LYS:HE3	1:C:230:GLU:OE1	1.87	0.75
1:B:33:PHE:N	1:B:41:ASN:HD21	1.83	0.75
1:C:229:GLN:CD	1:C:229:GLN:H	1.94	0.74
1:C:97:ALA:O	1:C:101:MET:HB2	1.87	0.74
1:C:10:ARG:HH12	1:C:424:ARG:HG2	1.52	0.73
1:A:155:ARG:HD3	1:A:431:VAL:O	1.88	0.73
1:A:34:GLN:HE21	1:A:63:ASP:HB3	1.54	0.73
1:C:404:GLU:CD	1:C:404:GLU:H	1.96	0.73
1:C:10:ARG:HH22	1:C:424:ARG:NH1	1.86	0.73
1:D:407:LEU:HD22	1:D:422:LEU:HD21	1.70	0.73
1:D:360:ILE:HG22	1:D:361:LEU:HD13	1.70	0.72
1:B:208:LEU:HD23	1:B:218:VAL:HG21	1.71	0.72
1:D:309:VAL:HG11	1:D:324:HIS:CD2	2.25	0.71
1:C:43:ALA:HB1	1:C:44:PRO:HD2	1.72	0.71
1:A:34:GLN:NE2	1:A:63:ASP:HB3	2.06	0.71
1:B:32:MET:HE2	1:B:105:PHE:HZ	1.56	0.70
1:C:34:GLN:HE21	1:C:63:ASP:HB3	1.55	0.70
1:C:218:VAL:HB	1:C:308:VAL:HG12	1.74	0.70
1:D:318:GLY:HA2	1:D:322:LEU:HD11	1.74	0.69
1:B:33:PHE:H	1:B:41:ASN:ND2	1.86	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:278:GLN:HB2	1:B:280:LYS:HD3	1.74	0.69
1:C:33:PHE:N	1:C:41:ASN:HD21	1.91	0.68
1:C:163:LEU:HD21	1:C:389:TRP:CD2	2.29	0.68
1:C:33:PHE:H	1:C:41:ASN:ND2	1.89	0.68
1:C:209:GLU:HG2	1:C:243:PRO:HD3	1.77	0.67
1:C:220:ILE:HG12	1:C:337:VAL:CG2	2.25	0.67
1:C:221:PRO:HB3	1:C:321:ILE:HG12	1.76	0.67
1:D:302:ARG:HG2	1:D:303:ALA:N	2.08	0.67
1:C:362:GLY:O	1:C:363:GLU:HG3	1.95	0.67
1:A:33:PHE:N	1:A:41:ASN:HD21	1.86	0.66
1:D:325:LEU:HD13	1:D:360:ILE:HD13	1.77	0.66
1:D:250:ASP:HA	1:D:291:VAL:HB	1.75	0.66
1:B:42:HIS:ND1	1:B:105:PHE:HB3	2.09	0.66
1:D:410:GLY:HA2	1:D:420:VAL:HG21	1.78	0.66
1:B:289:GLU:HG3	6:B:518:HOH:O	1.94	0.66
1:A:33:PHE:H	1:A:41:ASN:ND2	1.87	0.66
1:D:222:THR:HG22	1:D:339:VAL:CG2	2.26	0.66
1:C:247:ILE:HA	1:C:308:VAL:HG23	1.78	0.65
1:D:309:VAL:C	1:D:310:LEU:HD12	2.22	0.65
1:D:99:LYS:HE3	2:D:443:SO4:O2	1.97	0.65
1:A:86:THR:HG22	1:A:90:MET:HE2	1.79	0.64
1:C:322:LEU:HB3	1:C:361:LEU:HD11	1.78	0.64
1:C:235:LEU:HD13	1:C:247:ILE:HD13	1.78	0.64
1:D:401:GLY:HA3	1:D:406:LEU:HD11	1.79	0.64
1:C:83:THR:O	1:C:87:VAL:HG23	1.98	0.64
1:C:229:GLN:H	1:C:229:GLN:NE2	1.95	0.64
1:D:298:LYS:HA	1:D:301:ASN:HD22	1.62	0.64
1:D:200:THR:HG21	1:D:376:GLY:HA3	1.78	0.63
1:D:224:ALA:HB1	1:D:253:MET:HG2	1.80	0.63
1:D:325:LEU:O	1:D:329:LEU:HD13	1.97	0.63
1:D:239:GLY:HA2	1:D:242:LEU:HD12	1.80	0.63
1:A:362:GLY:HA2	2:A:443:SO4:O3	1.98	0.63
1:D:422:LEU:HD12	1:D:422:LEU:N	2.14	0.63
1:D:331:ASP:HB3	1:D:334:ASN:ND2	2.13	0.62
1:C:32:MET:HE1	1:C:101:MET:HE1	1.80	0.62
1:D:211:THR:HG21	1:D:335:ALA:CB	2.27	0.62
1:A:381:ALA:HB3	1:A:386:LEU:HD13	1.82	0.62
1:D:321:ILE:O	1:D:325:LEU:HG	2.00	0.62
1:C:358:VAL:O	1:C:365:VAL:HG22	2.00	0.62
1:A:321:ILE:O	1:A:325:LEU:HD13	2.00	0.62
1:B:87:VAL:HA	1:B:90:MET:HE3	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:253:MET:HA	1:D:256:ARG:HH21	1.65	0.62
1:B:165:ASN:HD22	1:B:165:ASN:C	2.07	0.61
1:D:209:GLU:HG3	1:D:243:PRO:HD3	1.81	0.61
1:D:97:ALA:O	1:D:101:MET:HB2	2.01	0.61
1:D:12:VAL:HG23	1:D:13:THR:HG23	1.83	0.60
1:A:32:MET:HE1	1:A:101:MET:HE1	1.84	0.60
1:C:1:MET:HG2	1:C:431:VAL:HG21	1.83	0.60
1:A:253:MET:HA	1:A:256:ARG:NH2	2.17	0.60
1:D:139:ALA:O	1:D:174:PRO:HG3	2.01	0.60
1:D:10:ARG:NH1	1:D:424:ARG:HG3	2.17	0.60
1:D:208:LEU:HD23	1:D:218:VAL:HG11	1.83	0.60
1:D:253:MET:HA	1:D:256:ARG:NH2	2.16	0.60
1:A:306:PRO:HD2	1:B:302:ARG:NH2	2.12	0.60
1:C:34:GLN:NE2	1:C:63:ASP:HB3	2.17	0.59
1:D:85:ALA:HB3	1:D:144:GLY:HA3	1.83	0.59
1:D:229:GLN:H	1:D:229:GLN:CD	2.10	0.59
1:A:182:LEU:HD11	1:A:397:VAL:HG23	1.84	0.59
1:B:269:SER:O	1:B:273:GLN:HG3	2.01	0.59
1:A:259:SER:O	1:A:262:PRO:HD2	2.03	0.59
1:C:87:VAL:HG13	1:C:118:LEU:HD13	1.84	0.59
1:D:33:PHE:H	1:D:41:ASN:ND2	1.90	0.59
1:C:209:GLU:HG2	1:C:243:PRO:CD	2.33	0.58
1:C:424:ARG:HD3	1:C:427:GLU:OE2	2.02	0.58
1:C:32:MET:HE3	1:C:62:LEU:HD22	1.86	0.58
1:B:289:GLU:CG	6:B:518:HOH:O	2.50	0.58
1:C:85:ALA:HB2	1:C:267:TYR:CD2	2.38	0.58
1:C:222:THR:HG22	1:C:339:VAL:HG21	1.86	0.58
1:C:170:VAL:HG13	1:C:171:LEU:HD13	1.86	0.58
1:D:45:PHE:HB3	1:D:47:PHE:CE1	2.39	0.58
1:B:235:LEU:HD13	1:B:247:ILE:HD13	1.86	0.57
1:B:359:ARG:HD2	1:B:364:GLU:OE2	2.04	0.57
1:C:32:MET:HE2	1:C:105:PHE:HZ	1.68	0.57
1:D:59:HIS:CD2	1:D:61:HIS:HB2	2.39	0.57
1:D:229:GLN:H	1:D:229:GLN:NE2	2.02	0.57
1:A:57:LEU:HG	1:A:65:VAL:HG12	1.85	0.57
1:D:98:LEU:HD21	1:D:108:PRO:CB	2.34	0.57
1:B:1:MET:HG3	1:B:21:ALA:HB2	1.85	0.57
1:C:86:THR:HG22	1:C:90:MET:CE	2.34	0.57
1:D:34:GLN:HE21	1:D:63:ASP:HB3	1.70	0.57
1:B:12:VAL:HG21	4:B:456:U5P:O2	2.04	0.57
1:C:217:LYS:HB2	1:C:334:ASN:OD1	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:LEU:HD21	1:A:389:TRP:CD2	2.40	0.57
1:C:98:LEU:HD11	1:C:108:PRO:CA	2.26	0.57
1:C:355:PRO:HB2	1:C:356:PRO:HD2	1.87	0.57
1:C:10:ARG:NH2	1:C:424:ARG:NH1	2.53	0.56
1:C:40:ARG:HD2	6:C:467:HOH:O	2.05	0.56
1:C:220:ILE:HG23	1:C:337:VAL:HG23	1.87	0.56
1:A:129:LEU:HD21	6:A:575:HOH:O	2.05	0.56
1:D:355:PRO:HB2	1:D:356:PRO:HD2	1.86	0.56
1:A:189:TYR:CZ	4:A:465:U5P:H5	2.40	0.56
1:C:163:LEU:HD21	1:C:389:TRP:CE3	2.40	0.56
1:C:225:VAL:O	1:C:257:VAL:HG11	2.06	0.56
1:A:284:ARG:HD3	1:A:288:LEU:HD23	1.88	0.56
1:A:195:ARG:HD2	1:A:199:GLU:OE2	2.05	0.56
1:C:88:LEU:HB3	1:C:260:LEU:HD21	1.88	0.56
1:D:11:GLU:CD	1:D:40:ARG:HH12	2.14	0.56
1:D:240:HIS:ND1	1:D:241:ARG:N	2.54	0.56
1:C:309:VAL:C	1:C:310:LEU:HD12	2.30	0.56
1:C:321:ILE:O	1:C:325:LEU:HD13	2.05	0.55
1:C:59:HIS:CD2	1:C:61:HIS:HB2	2.41	0.55
1:C:402:GLU:HB3	1:C:404:GLU:OE2	2.07	0.55
1:C:86:THR:HG22	1:C:90:MET:HE2	1.89	0.55
1:A:165:ASN:C	1:A:165:ASN:HD22	2.15	0.55
1:C:298:LYS:HD2	2:C:453:SO4:O3	2.06	0.55
1:D:170:VAL:HG12	1:D:171:LEU:HD12	1.87	0.55
1:D:396:VAL:HG12	1:D:398:LEU:HD12	1.89	0.55
1:C:331:ASP:HB3	1:C:334:ASN:ND2	2.21	0.55
1:C:360:ILE:CG2	1:C:361:LEU:HD13	2.34	0.54
1:B:27:LEU:HD13	1:B:29:ASP:O	2.07	0.54
1:D:360:ILE:O	1:D:361:LEU:HB2	2.07	0.54
1:A:370:SER:HA	2:A:451:SO4:O2	2.07	0.54
1:B:202:ARG:O	1:B:206:GLU:HG3	2.07	0.54
1:C:401:GLY:HA3	1:C:406:LEU:HD11	1.89	0.54
1:D:284:ARG:HD3	1:D:288:LEU:CD2	2.33	0.54
1:C:236:TYR:HD1	1:C:285:PRO:HA	1.73	0.54
1:B:208:LEU:HD21	1:B:218:VAL:HG11	1.90	0.54
1:D:229:GLN:HG3	1:D:261:TYR:CZ	2.42	0.54
1:A:168:LYS:HE3	1:A:230:GLU:CD	2.33	0.53
1:C:182:LEU:HD11	1:C:397:VAL:HG23	1.89	0.53
1:D:224:ALA:CB	1:D:253:MET:HG2	2.38	0.53
1:D:236:TYR:OH	1:D:280:LYS:HE2	2.08	0.53
1:A:359:ARG:NH1	2:A:443:SO4:O3	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:69:PRO:O	1:D:110:ASP:HB3	2.09	0.53
1:D:225:VAL:O	1:D:257:VAL:HG11	2.08	0.53
1:C:214:GLN:NE2	1:C:333:ARG:HA	2.24	0.53
1:D:2:ARG:HG3	1:D:430:PRO:HA	1.90	0.53
1:D:1:MET:HG3	1:D:21:ALA:CB	2.35	0.53
1:D:360:ILE:CG2	1:D:361:LEU:HD13	2.38	0.53
1:C:270:GLU:HA	1:C:270:GLU:OE1	2.09	0.52
1:B:226:GLU:HG2	6:B:538:HOH:O	2.08	0.52
1:A:344:GLN:NE2	1:A:344:GLN:HA	2.24	0.52
1:C:236:TYR:CD1	1:C:285:PRO:HA	2.45	0.52
1:D:165:ASN:C	1:D:165:ASN:HD22	2.16	0.52
1:A:98:LEU:HD11	1:A:108:PRO:HA	1.91	0.52
1:B:165:ASN:C	1:B:165:ASN:ND2	2.66	0.52
1:A:263:ARG:HG2	1:D:277:LEU:HD21	1.91	0.52
1:C:57:LEU:HG	1:C:65:VAL:HG12	1.92	0.52
1:B:274:ALA:O	1:B:278:GLN:HG2	2.10	0.52
1:B:411:LYS:HB2	3:B:455:FLC:OHB	2.10	0.52
1:C:5:PRO:HG2	1:C:423:ALA:HB1	1.91	0.52
1:B:42:HIS:CE1	1:B:105:PHE:HB3	2.45	0.52
1:C:28:LEU:O	1:C:29:ASP:HB2	2.09	0.52
1:D:214:GLN:HE21	1:D:333:ARG:HA	1.75	0.52
1:D:68:LEU:N	1:D:69:PRO:HD2	2.25	0.51
1:D:183:VAL:HG21	1:D:393:GLU:HG3	1.92	0.51
1:B:102:ASP:HA	6:B:524:HOH:O	2.09	0.51
1:C:294:THR:O	1:C:298:LYS:HG2	2.11	0.51
1:B:90:MET:HE1	1:B:118:LEU:HD22	1.92	0.51
1:D:101:MET:HE3	1:D:101:MET:HA	1.92	0.51
1:D:155:ARG:HD3	1:D:431:VAL:O	2.11	0.51
1:A:11:GLU:OE2	1:A:40:ARG:NH1	2.42	0.51
1:C:155:ARG:HD3	1:C:431:VAL:O	2.11	0.51
1:C:292:GLU:HB2	2:C:433:SO4:O2	2.11	0.51
1:A:297:SER:OG	1:A:320:ARG:HG2	2.11	0.51
1:C:84:ARG:NH1	2:C:451:SO4:O3	2.36	0.51
1:A:32:MET:HA	1:A:67:ARG:HG3	1.93	0.50
1:D:336:LEU:C	1:D:336:LEU:HD23	2.35	0.50
1:B:36:LYS:C	1:B:37:GLU:HG2	2.37	0.50
1:C:59:HIS:HD2	1:C:61:HIS:H	1.58	0.50
1:C:409:LEU:HD22	1:C:413:LEU:HG	1.92	0.50
1:D:214:GLN:NE2	1:D:333:ARG:HA	2.27	0.50
1:D:218:VAL:O	1:D:220:ILE:HG13	2.11	0.50
1:D:319:GLY:C	1:D:321:ILE:H	2.20	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:LEU:HD23	1:A:150:ALA:HA	1.92	0.50
1:B:208:LEU:CD2	1:B:218:VAL:HG11	2.41	0.50
1:C:247:ILE:HG12	1:C:308:VAL:CG2	2.41	0.50
1:B:211:THR:HA	1:B:214:GLN:HG2	1.93	0.50
1:C:9:ALA:O	1:C:11:GLU:HG2	2.11	0.50
1:C:170:VAL:HG13	1:C:171:LEU:CD1	2.41	0.50
1:C:170:VAL:HG22	1:C:171:LEU:HD12	1.93	0.50
1:D:208:LEU:CD2	1:D:218:VAL:HG11	2.41	0.50
1:A:344:GLN:HE21	1:A:344:GLN:CA	2.25	0.50
1:B:12:VAL:HG23	1:B:13:THR:HG23	1.93	0.50
1:C:10:ARG:HH22	1:C:424:ARG:HH11	1.59	0.50
1:D:86:THR:HG22	1:D:90:MET:CE	2.39	0.50
1:D:42:HIS:CE1	1:D:105:PHE:HB3	2.47	0.50
1:D:228:ALA:HB3	1:D:229:GLN:NE2	2.27	0.50
1:A:359:ARG:NH2	2:A:444:SO4:O2	2.45	0.49
1:C:381:ALA:HB3	1:C:386:LEU:HD13	1.94	0.49
1:D:298:LYS:HA	1:D:301:ASN:ND2	2.25	0.49
1:D:226:GLU:N	6:D:476:HOH:O	2.44	0.49
1:D:299:ALA:HA	1:D:302:ARG:HD2	1.94	0.49
1:B:208:LEU:CD2	1:B:218:VAL:HG21	2.41	0.49
1:A:59:HIS:CD2	1:A:61:HIS:HB2	2.47	0.49
1:B:34:GLN:HA	1:B:38:GLU:HG3	1.94	0.49
1:D:219:LEU:HD21	1:D:324:HIS:O	2.12	0.49
1:A:396:VAL:O	1:A:420:VAL:HA	2.11	0.49
1:D:33:PHE:N	1:D:41:ASN:HD21	1.93	0.49
1:D:326:LYS:HD2	1:D:361:LEU:HD22	1.93	0.49
1:A:306:PRO:CD	1:B:302:ARG:HH22	2.15	0.49
1:C:250:ASP:HA	1:C:291:VAL:HB	1.94	0.49
1:D:322:LEU:HA	1:D:325:LEU:HD11	1.94	0.49
1:C:128:ARG:NH2	6:C:471:HOH:O	2.45	0.49
1:B:34:GLN:HE21	1:B:63:ASP:HB3	1.78	0.49
1:B:163:LEU:HD21	1:B:389:TRP:CE3	2.48	0.49
1:D:65:VAL:CG1	1:D:94:LEU:HD11	2.43	0.49
1:D:240:HIS:CE1	1:D:241:ARG:HB3	2.48	0.49
1:B:11:GLU:OE1	1:B:40:ARG:NH1	2.45	0.48
1:A:389:TRP:HE3	1:A:390:LEU:HD13	1.78	0.48
1:D:359:ARG:HD2	1:D:359:ARG:C	2.38	0.48
1:D:12:VAL:HG12	1:D:401:GLY:HA2	1.95	0.48
1:D:203:GLU:O	1:D:207:ILE:HG13	2.13	0.48
1:D:239:GLY:HA2	1:D:242:LEU:CD1	2.43	0.48
1:D:329:LEU:HA	1:D:369:ALA:CB	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:366:PRO:HG2	2:C:454:SO4:O2	2.14	0.48
1:D:197:TYR:O	1:D:201:VAL:HG23	2.14	0.48
1:D:402:GLU:O	1:D:406:LEU:HD13	2.14	0.48
1:C:312:GLY:HA2	1:C:313:SER:C	2.38	0.48
1:D:221:PRO:HD2	1:D:337:VAL:O	2.13	0.48
1:C:84:ARG:HB3	1:C:267:TYR:OH	2.13	0.48
1:C:323:HIS:ND1	2:C:453:SO4:O2	2.47	0.48
1:C:365:VAL:HG23	1:C:365:VAL:O	2.12	0.48
1:D:384:ASP:OD2	1:D:384:ASP:N	2.46	0.48
1:C:239:GLY:C	1:C:241:ARG:H	2.20	0.48
1:D:62:LEU:HD13	1:D:93:VAL:HG12	1.95	0.48
1:C:57:LEU:HD23	1:C:90:MET:CE	2.44	0.48
1:C:113:GLU:OE2	1:C:117:HIS:HE1	1.96	0.48
1:C:325:LEU:HG	1:C:329:LEU:HD11	1.96	0.48
1:D:280:LYS:O	1:D:282:PRO:HD3	2.14	0.48
1:A:217:LYS:HE2	1:A:307:MET:HE3	1.96	0.48
1:D:3:ILE:HG12	1:D:184:LEU:HD22	1.95	0.48
1:D:34:GLN:NE2	1:D:63:ASP:HB3	2.29	0.48
1:A:43:ALA:HB1	1:A:44:PRO:HD2	1.95	0.48
1:B:168:LYS:HG2	1:B:197:TYR:CE2	2.48	0.48
1:C:142:LEU:HG	1:C:143:PRO:HD2	1.96	0.48
1:B:11:GLU:OE2	1:B:37:GLU:HG3	2.13	0.47
1:D:227:ARG:HG2	1:D:227:ARG:HH11	1.79	0.47
1:B:168:LYS:HE2	1:B:230:GLU:OE1	2.14	0.47
1:C:247:ILE:HG12	1:C:308:VAL:HG21	1.96	0.47
1:D:211:THR:OG1	1:D:218:VAL:HG22	2.15	0.47
1:B:142:LEU:HD22	1:B:226:GLU:HB2	1.96	0.47
1:A:163:LEU:HD21	1:A:389:TRP:CE3	2.48	0.47
1:C:226:GLU:HG2	1:C:261:TYR:OH	2.13	0.47
1:A:389:TRP:CE3	1:A:390:LEU:HD13	2.50	0.47
1:B:221:PRO:HB3	1:B:321:ILE:HG12	1.96	0.47
1:B:344:GLN:CD	1:B:344:GLN:H	2.21	0.47
1:D:329:LEU:HA	1:D:369:ALA:HB2	1.97	0.47
1:D:360:ILE:HD12	1:D:365:VAL:HG11	1.97	0.47
1:D:399:VAL:HG12	1:D:423:ALA:HB3	1.97	0.47
1:C:211:THR:HG21	1:C:335:ALA:CB	2.45	0.47
1:C:262:PRO:HG3	1:C:281:ASN:ND2	2.30	0.47
1:D:238:HIS:C	1:D:240:HIS:H	2.23	0.47
1:D:359:ARG:HD3	1:D:362:GLY:HA2	1.97	0.47
1:C:253:MET:O	1:C:257:VAL:HG23	2.15	0.46
1:D:202:ARG:O	1:D:206:GLU:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:231:ILE:O	1:D:235:LEU:HD13	2.14	0.46
1:A:11:GLU:CD	1:A:40:ARG:HH12	2.23	0.46
1:A:191:ASP:CG	1:A:405:LYS:HD2	2.40	0.46
1:B:420:VAL:HG22	1:B:421:SER:N	2.30	0.46
1:B:404:GLU:OE2	1:B:404:GLU:HA	2.16	0.46
1:D:65:VAL:HG11	1:D:94:LEU:HD11	1.97	0.46
1:D:347:LEU:HD12	1:D:347:LEU:H	1.81	0.46
1:B:68:LEU:N	1:B:69:PRO:HD2	2.30	0.46
1:A:45:PHE:HB3	1:A:47:PHE:CE1	2.51	0.46
1:A:85:ALA:HB3	1:A:144:GLY:HA3	1.97	0.46
1:A:401:GLY:HA3	1:A:406:LEU:HD13	1.97	0.46
1:D:204:PHE:O	1:D:208:LEU:HG	2.15	0.46
1:D:318:GLY:HA2	1:D:322:LEU:CD1	2.43	0.46
1:A:153:GLU:CG	2:A:433:SO4:O1	2.50	0.46
1:C:57:LEU:HD21	1:C:68:LEU:HD22	1.98	0.46
1:C:192:ARG:HA	1:C:383:GLN:NE2	2.31	0.46
1:C:214:GLN:HE21	1:C:333:ARG:HA	1.81	0.46
1:D:347:LEU:HD12	1:D:347:LEU:N	2.30	0.46
1:B:98:LEU:HD11	1:B:108:PRO:HA	1.97	0.46
1:C:31:GLY:HA3	1:C:63:ASP:C	2.41	0.46
1:C:277:LEU:C	1:C:279:GLY:H	2.24	0.46
1:D:28:LEU:O	1:D:29:ASP:HB2	2.14	0.46
1:C:85:ALA:HB2	1:C:267:TYR:CE2	2.51	0.45
1:D:295:GLU:H	1:D:295:GLU:CD	2.22	0.45
1:B:217:LYS:HG2	1:B:307:MET:HG2	1.97	0.45
1:B:411:LYS:HD2	3:B:455:FLC:OHB	2.15	0.45
1:B:348:GLY:O	1:B:352:ILE:HG13	2.16	0.45
1:C:177:PRO:HD3	1:C:389:TRP:CE2	2.51	0.45
1:A:289:GLU:HG3	6:A:502:HOH:O	2.17	0.45
1:D:217:LYS:HG2	1:D:307:MET:HG2	1.99	0.45
1:D:234:VAL:O	1:D:238:HIS:HB2	2.17	0.45
1:D:347:LEU:O	1:D:351:ILE:HG13	2.16	0.45
1:B:163:LEU:HD21	1:B:389:TRP:CD2	2.51	0.45
1:B:140:GLY:O	1:B:164:GLY:HA3	2.17	0.45
1:C:214:GLN:O	1:C:333:ARG:HD2	2.17	0.45
1:B:276:PHE:HA	1:B:280:LYS:O	2.17	0.45
1:C:170:VAL:C	1:C:171:LEU:HD12	2.42	0.45
1:A:31:GLY:HA3	1:A:63:ASP:O	2.16	0.45
1:B:178:PRO:HB3	1:C:126:TRP:CE3	2.52	0.45
1:C:140:GLY:O	1:C:164:GLY:HA3	2.17	0.45
1:C:172:PRO:HD3	1:C:268:PHE:C	2.42	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:399:VAL:HG12	1:C:423:ALA:HB3	1.98	0.45
1:C:420:VAL:HG22	1:C:421:SER:N	2.31	0.45
1:A:192:ARG:NH2	2:A:463:SO4:O2	2.47	0.45
1:A:407:LEU:HD13	1:A:422:LEU:CD2	2.47	0.45
1:B:182:LEU:HD11	1:B:397:VAL:HG23	1.99	0.45
1:C:143:PRO:HD3	1:C:226:GLU:HG3	1.99	0.45
1:C:209:GLU:O	1:C:210:LYS:C	2.59	0.45
1:C:229:GLN:HG3	1:C:261:TYR:CZ	2.53	0.45
1:C:267:TYR:HA	6:C:500:HOH:O	2.16	0.45
1:D:326:LYS:O	1:D:326:LYS:HG2	2.17	0.45
1:D:397:VAL:HG21	1:D:429:VAL:HG11	1.98	0.45
1:A:235:LEU:HD13	1:A:247:ILE:HD13	1.99	0.44
1:B:428:GLY:H	4:B:457:U5P:C5	2.30	0.44
1:C:404:GLU:CD	1:C:404:GLU:N	2.72	0.44
1:D:386:LEU:O	1:D:390:LEU:HD22	2.17	0.44
1:A:262:PRO:O	1:A:265:VAL:HG23	2.18	0.44
1:A:318:GLY:HA2	1:A:322:LEU:CD1	2.47	0.44
1:A:341:TYR:HB3	4:A:465:U5P:H5'2	2.00	0.44
1:C:211:THR:HG21	1:C:335:ALA:HB3	1.98	0.44
1:B:318:GLY:HA2	1:B:322:LEU:CD1	2.48	0.44
1:C:331:ASP:HB3	1:C:334:ASN:HD22	1.83	0.44
1:B:37:GLU:O	1:B:39:ALA:N	2.50	0.44
1:D:404:GLU:H	1:D:404:GLU:CD	2.25	0.44
1:C:1:MET:HG2	1:C:431:VAL:CG2	2.47	0.44
1:C:302:ARG:HG2	1:C:302:ARG:HH21	1.83	0.44
1:B:259:SER:O	1:B:262:PRO:HD2	2.17	0.44
1:C:294:THR:O	1:C:297:SER:HB3	2.17	0.44
1:A:102:ASP:CG	1:A:103:GLU:H	2.26	0.44
1:B:91:GLU:O	1:B:95:GLU:HG2	2.18	0.44
1:C:251:SER:HB3	1:C:254:ALA:HB3	2.00	0.44
1:D:235:LEU:O	1:D:239:GLY:HA3	2.17	0.44
1:A:165:ASN:C	1:A:165:ASN:ND2	2.74	0.44
1:B:191:ASP:CG	1:B:192:ARG:HG3	2.43	0.44
1:D:220:ILE:HG22	1:D:222:THR:HG23	2.00	0.44
1:D:359:ARG:HD3	1:D:362:GLY:CA	2.48	0.44
1:A:102:ASP:HB2	6:A:580:HOH:O	2.17	0.43
1:D:422:LEU:N	1:D:422:LEU:CD1	2.80	0.43
1:A:85:ALA:HB2	1:A:267:TYR:CD2	2.53	0.43
1:B:344:GLN:H	1:B:344:GLN:NE2	2.17	0.43
1:D:3:ILE:HG23	1:D:3:ILE:O	2.18	0.43
1:D:233:TYR:CE1	1:D:282:PRO:HB2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:ARG:HG2	6:A:501:HOH:O	2.17	0.43
1:A:344:GLN:HA	1:A:344:GLN:HE21	1.79	0.43
1:C:10:ARG:NH2	1:C:424:ARG:HH12	2.15	0.43
1:D:158:VAL:HG23	1:D:180:ALA:HB2	2.01	0.43
1:D:252:PRO:O	1:D:256:ARG:HG3	2.18	0.43
1:A:84:ARG:HE	1:A:84:ARG:HB2	1.40	0.43
1:A:289:GLU:CG	6:A:502:HOH:O	2.67	0.43
1:C:362:GLY:C	1:C:363:GLU:HG3	2.43	0.43
1:B:101:MET:HG2	1:B:104:PRO:CB	2.41	0.43
1:B:113:GLU:OE2	1:B:117:HIS:HE1	2.01	0.43
1:D:235:LEU:HD23	1:D:247:ILE:HG21	2.00	0.43
1:D:325:LEU:HD12	1:D:326:LYS:N	2.34	0.43
1:B:70:LYS:NZ	1:B:74:GLU:OE1	2.43	0.43
1:D:195:ARG:O	1:D:196:PRO:C	2.61	0.43
1:A:58:THR:O	1:A:145:SER:HA	2.18	0.43
1:A:212:LEU:CB	1:A:243:PRO:HG2	2.49	0.43
1:C:229:GLN:HA	1:C:232:LEU:HD12	2.01	0.43
1:D:250:ASP:HB3	1:D:311:ALA:HB2	1.99	0.43
1:B:270:GLU:OE1	1:C:266:ARG:NH2	2.51	0.43
1:D:356:PRO:HG2	1:D:357:ALA:H	1.84	0.43
1:D:389:TRP:HE3	1:D:390:LEU:HD13	1.84	0.43
1:A:61:HIS:CD2	1:A:142:LEU:HD11	2.53	0.43
1:B:396:VAL:O	1:B:420:VAL:HA	2.19	0.43
1:D:216:GLY:HA3	1:D:333:ARG:O	2.19	0.43
1:D:316:LEU:HD22	1:D:325:LEU:HD21	2.01	0.43
1:A:98:LEU:HD11	1:A:108:PRO:CA	2.49	0.43
1:A:197:TYR:O	1:A:200:THR:HB	2.19	0.43
1:A:203:GLU:O	1:A:207:ILE:HG13	2.19	0.43
1:C:386:LEU:O	1:C:390:LEU:HD23	2.18	0.43
1:D:14:GLY:O	1:D:15:SER:C	2.61	0.43
1:A:87:VAL:HA	1:A:90:MET:HE3	2.01	0.42
1:B:81:TYR:HA	1:B:119:ARG:O	2.19	0.42
1:C:272:VAL:O	1:C:276:PHE:HD2	2.00	0.42
1:A:31:GLY:HA3	1:A:63:ASP:C	2.44	0.42
1:B:33:PHE:HB3	1:B:37:GLU:HB2	2.00	0.42
1:D:396:VAL:HG12	1:D:398:LEU:CD1	2.49	0.42
1:A:80:VAL:HB	1:A:118:LEU:HD23	2.01	0.42
1:D:5:PRO:HG2	1:D:423:ALA:HB1	2.00	0.42
1:D:309:VAL:HG12	1:D:310:LEU:N	2.32	0.42
1:D:401:GLY:CA	1:D:406:LEU:HD11	2.45	0.42
1:A:409:LEU:HD22	1:A:413:LEU:HG	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:HIS:CE1	1:A:162:ASP:OD1	2.73	0.42
1:C:12:VAL:HG12	1:C:401:GLY:HA2	2.02	0.42
1:D:32:MET:HB2	1:D:41:ASN:OD1	2.20	0.42
1:A:77:ARG:HD2	2:A:459:SO4:O4	2.20	0.42
1:B:126:TRP:CE3	1:C:178:PRO:HB3	2.54	0.42
1:B:318:GLY:HA2	1:B:322:LEU:HD11	2.02	0.42
1:B:347:LEU:HD11	1:B:358:VAL:HG11	2.01	0.42
1:D:165:ASN:C	1:D:165:ASN:ND2	2.76	0.42
1:D:220:ILE:HG22	1:D:222:THR:CG2	2.49	0.42
1:A:46:GLY:HA2	3:A:464:FLC:HG2	2.00	0.42
1:D:65:VAL:O	1:D:65:VAL:HG12	2.19	0.42
1:A:212:LEU:HB2	1:A:243:PRO:HG2	2.01	0.42
1:A:365:VAL:HA	1:A:366:PRO:HD3	1.87	0.42
1:C:59:HIS:NE2	1:C:61:HIS:HB2	2.35	0.42
1:C:386:LEU:O	1:C:390:LEU:CD2	2.68	0.42
1:D:82:ALA:O	1:D:120:PRO:HA	2.20	0.42
1:A:321:ILE:HG23	1:A:322:LEU:N	2.34	0.42
1:C:326:LYS:HD2	1:C:361:LEU:HD22	2.02	0.42
1:A:140:GLY:O	1:A:164:GLY:HA3	2.19	0.41
1:B:4:VAL:HA	1:B:5:PRO:HD3	1.85	0.41
1:C:187:GLY:O	1:C:188:THR:C	2.63	0.41
1:B:203:GLU:O	1:B:207:ILE:HG13	2.20	0.41
1:D:58:THR:O	1:D:145:SER:HA	2.19	0.41
1:D:297:SER:CB	1:D:320:ARG:HD2	2.50	0.41
1:D:302:ARG:HG2	1:D:303:ALA:H	1.81	0.41
1:A:407:LEU:HD13	1:A:422:LEU:HD21	2.02	0.41
1:C:10:ARG:CD	1:C:422:LEU:HD23	2.51	0.41
1:C:142:LEU:O	1:C:143:PRO:C	2.58	0.41
1:A:354:ARG:N	1:A:355:PRO:CD	2.83	0.41
1:C:203:GLU:O	1:C:207:ILE:HG13	2.21	0.41
1:C:93:VAL:HA	1:C:253:MET:HE1	2.01	0.41
1:A:57:LEU:HD23	1:A:90:MET:CE	2.51	0.41
1:D:170:VAL:HG12	1:D:171:LEU:CD1	2.51	0.41
1:C:350:GLU:O	1:C:353:ALA:HB3	2.21	0.41
1:D:347:LEU:H	1:D:347:LEU:CD1	2.34	0.41
1:A:91:GLU:O	1:A:95:GLU:HG2	2.21	0.41
1:A:318:GLY:HA2	1:A:322:LEU:HD11	2.02	0.41
1:B:10:ARG:HH11	1:B:10:ARG:HG2	1.86	0.41
1:B:55:VAL:HG12	1:B:57:LEU:HD13	2.02	0.41
1:B:168:LYS:HE2	1:B:230:GLU:CD	2.46	0.41
1:B:365:VAL:HA	1:B:366:PRO:HD3	1.80	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:49:PRO:HB3	1:C:71:LEU:HD12	2.02	0.41
1:C:98:LEU:HD12	1:C:111:VAL:HG21	2.03	0.41
1:C:217:LYS:HE2	1:C:307:MET:HE3	2.03	0.41
1:C:239:GLY:C	1:C:241:ARG:N	2.79	0.41
1:D:98:LEU:CD2	1:D:108:PRO:HA	2.51	0.41
1:D:397:VAL:HA	1:D:421:SER:O	2.20	0.41
1:D:409:LEU:HD22	1:D:413:LEU:HD11	2.02	0.41
1:C:247:ILE:HG23	1:C:308:VAL:HG23	2.04	0.41
1:C:318:GLY:HA2	1:C:322:LEU:HD11	2.01	0.41
1:C:326:LYS:HE3	1:C:363:GLU:OE1	2.21	0.41
1:D:33:PHE:HD2	1:D:41:ASN:ND2	2.19	0.41
1:A:98:LEU:C	1:A:98:LEU:HD23	2.46	0.40
1:A:420:VAL:HG22	1:A:421:SER:N	2.36	0.40
1:C:91:GLU:O	1:C:95:GLU:HG2	2.21	0.40
1:C:227:ARG:HG2	1:C:227:ARG:HH21	1.86	0.40
1:D:86:THR:CG2	1:D:90:MET:HE2	2.42	0.40
1:D:409:LEU:O	1:D:413:LEU:HG	2.21	0.40
1:A:126:TRP:CE3	1:A:135:ALA:HB2	2.56	0.40
1:A:424:ARG:HE	1:A:427:GLU:CD	2.30	0.40
1:C:360:ILE:HD12	1:C:365:VAL:HG21	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	429/431 (100%)	411 (96%)	17 (4%)	1 (0%)	43	52
1	B	429/431 (100%)	413 (96%)	15 (4%)	1 (0%)	43	52
1	C	429/431 (100%)	395 (92%)	32 (8%)	2 (0%)	24	27
1	D	429/431 (100%)	378 (88%)	44 (10%)	7 (2%)	7	6
All	All	1716/1724 (100%)	1597 (93%)	108 (6%)	11 (1%)	21	24

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	38	GLU
1	D	102	ASP
1	A	102	ASP
1	C	102	ASP
1	C	198	ARG
1	D	166	ARG
1	D	244	ARG
1	D	38	GLU
1	D	316	LEU
1	D	15	SER
1	D	225	VAL

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	342/342 (100%)	321 (94%)	21 (6%)	17	20
1	B	342/342 (100%)	320 (94%)	22 (6%)	16	18
1	C	342/342 (100%)	326 (95%)	16 (5%)	23	30
1	D	342/342 (100%)	327 (96%)	15 (4%)	25	33
All	All	1368/1368 (100%)	1294 (95%)	74 (5%)	20	24

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LEU
1	A	57	LEU
1	A	84	ARG
1	A	163	LEU
1	A	165	ASN
1	A	171	LEU
1	A	186	GLU
1	A	192	ARG
1	A	219	LEU

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Mol	Chain	Res	Type
1	A	241	ARG
1	A	264	LEU
1	A	320	ARG
1	A	329	LEU
1	A	336	LEU
1	A	344	GLN
1	A	386	LEU
1	A	390	LEU
1	A	404	GLU
1	A	406	LEU
1	A	407	LEU
1	A	409	LEU
1	B	27	LEU
1	B	38	GLU
1	B	57	LEU
1	B	65	VAL
1	B	94	LEU
1	B	98	LEU
1	B	101	MET
1	B	102	ASP
1	B	163	LEU
1	B	165	ASN
1	B	186	GLU
1	B	192	ARG
1	B	219	LEU
1	B	264	LEU
1	B	325	LEU
1	B	329	LEU
1	B	365	VAL
1	B	386	LEU
1	B	390	LEU
1	B	406	LEU
1	B	407	LEU
1	B	409	LEU
1	C	27	LEU
1	C	28	LEU
1	C	57	LEU
1	C	94	LEU
1	C	98	LEU
1	C	112	GLU
1	C	163	LEU
1	C	186	GLU

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Mol	Chain	Res	Type
1	C	229	GLN
1	C	336	LEU
1	C	344	GLN
1	C	386	LEU
1	C	398	LEU
1	C	407	LEU
1	C	409	LEU
1	C	415	LEU
1	D	2	ARG
1	D	27	LEU
1	D	57	LEU
1	D	96	ASP
1	D	165	ASN
1	D	186	GLU
1	D	195	ARG
1	D	229	GLN
1	D	253	MET
1	D	264	LEU
1	D	293	HIS
1	D	302	ARG
1	D	344	GLN
1	D	390	LEU
1	D	409	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	A	41	ASN
1	A	59	HIS
1	A	61	HIS
1	A	165	ASN
1	A	344	GLN
1	A	383	GLN
1	B	34	GLN
1	B	41	ASN
1	B	59	HIS
1	B	61	HIS
1	B	165	ASN
1	B	275	HIS
1	B	344	GLN
1	B	383	GLN

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Mol	Chain	Res	Type
1	B	391	GLN
1	C	34	GLN
1	C	41	ASN
1	C	165	ASN
1	C	214	GLN
1	C	229	GLN
1	C	275	HIS
1	C	301	ASN
1	C	383	GLN
1	D	34	GLN
1	D	41	ASN
1	D	64	HIS
1	D	165	ASN
1	D	194	HIS
1	D	214	GLN
1	D	229	GLN
1	D	372	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 113 ligands modelled in this entry, 8 are monoatomic - leaving 105 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	461	-	4,4,4	1.08	0	6,6,6	0.57	0
2	SO4	A	432	-	4,4,4	1.08	0	6,6,6	0.58	0
2	SO4	D	439	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	B	443	-	4,4,4	1.07	0	6,6,6	0.58	0
4	U5P	A	465	-	22,22,22	1.94	5 (22%)	32,33,33	1.77	6 (18%)
2	SO4	C	438	-	4,4,4	1.06	0	6,6,6	0.58	0
4	U5P	A	466	-	22,22,22	2.04	5 (22%)	32,33,33	1.73	6 (18%)
2	SO4	A	439	-	4,4,4	1.05	0	6,6,6	0.59	0
2	SO4	B	451	-	4,4,4	1.08	0	6,6,6	0.57	0
2	SO4	B	435	-	4,4,4	1.06	0	6,6,6	0.58	0
2	SO4	C	439	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	A	452	-	4,4,4	1.08	0	6,6,6	0.57	0
2	SO4	C	432	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	A	448	-	4,4,4	1.09	0	6,6,6	0.57	0
2	SO4	C	449	-	4,4,4	1.06	0	6,6,6	0.58	0
2	SO4	D	443	-	4,4,4	1.05	0	6,6,6	0.57	0
2	SO4	B	439	-	4,4,4	1.08	0	6,6,6	0.56	0
4	U5P	D	448	-	22,22,22	1.99	5 (22%)	32,33,33	1.71	5 (15%)
4	U5P	A	467	-	22,22,22	1.99	5 (22%)	32,33,33	1.74	5 (15%)
2	SO4	C	442	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	D	441	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	A	438	-	4,4,4	1.08	0	6,6,6	0.56	0
2	SO4	C	455	-	4,4,4	1.05	0	6,6,6	0.58	0
2	SO4	A	433	-	4,4,4	1.06	0	6,6,6	0.57	0
3	FLC	B	455	-	12,12,12	1.64	5 (41%)	17,17,17	1.28	1 (5%)
2	SO4	B	452	-	4,4,4	1.05	0	6,6,6	0.60	0
2	SO4	A	458	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	A	454	-	4,4,4	1.04	0	6,6,6	0.56	0
2	SO4	A	455	-	4,4,4	1.07	0	6,6,6	0.58	0
2	SO4	B	433	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	D	446	-	4,4,4	1.07	0	6,6,6	0.57	0
4	U5P	C	457	-	22,22,22	1.93	5 (22%)	32,33,33	1.72	6 (18%)
2	SO4	A	449	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	B	442	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	D	432	-	4,4,4	1.06	0	6,6,6	0.58	0
2	SO4	B	438	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	D	436	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	D	435	-	4,4,4	1.07	0	6,6,6	0.58	0
2	SO4	B	453	-	4,4,4	1.08	0	6,6,6	0.57	0
4	U5P	B	456	-	22,22,22	1.92	4 (18%)	32,33,33	1.70	6 (18%)
2	SO4	A	446	-	4,4,4	1.07	0	6,6,6	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	U5P	B	457	-	22,22,22	2.11	5 (22%)	32,33,33	1.69	6 (18%)
2	SO4	A	462	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	C	446	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	C	441	-	4,4,4	1.06	0	6,6,6	0.58	0
3	FLC	A	464	-	12,12,12	1.55	5 (41%)	17,17,17	1.33	1 (5%)
2	SO4	C	453	-	4,4,4	1.07	0	6,6,6	0.58	0
2	SO4	A	453	-	4,4,4	1.08	0	6,6,6	0.59	0
2	SO4	B	446	-	4,4,4	1.05	0	6,6,6	0.59	0
2	SO4	D	444	-	4,4,4	1.06	0	6,6,6	0.59	0
2	SO4	A	441	-	4,4,4	1.07	0	6,6,6	0.58	0
2	SO4	C	434	-	4,4,4	1.05	0	6,6,6	0.58	0
2	SO4	A	442	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	D	438	-	4,4,4	1.05	0	6,6,6	0.58	0
2	SO4	A	437	-	4,4,4	1.08	0	6,6,6	0.56	0
2	SO4	D	434	-	4,4,4	1.09	0	6,6,6	0.57	0
2	SO4	B	454	-	4,4,4	1.08	0	6,6,6	0.57	0
2	SO4	A	450	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	B	440	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	D	440	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	A	445	-	4,4,4	1.07	0	6,6,6	0.58	0
2	SO4	B	437	-	4,4,4	1.06	0	6,6,6	0.58	0
2	SO4	B	436	-	4,4,4	1.02	0	6,6,6	0.61	0
2	SO4	B	441	-	4,4,4	1.07	0	6,6,6	0.58	0
2	SO4	A	451	-	4,4,4	1.04	0	6,6,6	0.58	0
2	SO4	C	440	-	4,4,4	1.06	0	6,6,6	0.58	0
2	SO4	D	437	-	4,4,4	1.06	0	6,6,6	0.56	0
2	SO4	B	450	-	4,4,4	1.08	0	6,6,6	0.57	0
2	SO4	C	452	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	C	445	-	4,4,4	1.07	0	6,6,6	0.59	0
2	SO4	C	451	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	A	460	-	4,4,4	1.10	0	6,6,6	0.57	0
2	SO4	B	447	-	4,4,4	1.07	0	6,6,6	0.58	0
2	SO4	A	434	-	4,4,4	1.08	0	6,6,6	0.57	0
2	SO4	C	433	-	4,4,4	1.08	0	6,6,6	0.58	0
2	SO4	B	448	-	4,4,4	1.08	0	6,6,6	0.56	0
2	SO4	A	435	-	4,4,4	1.09	0	6,6,6	0.57	0
2	SO4	A	440	-	4,4,4	1.10	0	6,6,6	0.59	0
2	SO4	A	447	-	4,4,4	1.08	0	6,6,6	0.57	0
2	SO4	C	450	-	4,4,4	1.08	0	6,6,6	0.57	0
2	SO4	C	443	-	4,4,4	1.08	0	6,6,6	0.56	0
2	SO4	C	447	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	C	454	-	4,4,4	1.07	0	6,6,6	0.58	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	B	445	-	4,4,4	1.10	0	6,6,6	0.53	0
2	SO4	C	437	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	A	436	-	4,4,4	1.08	0	6,6,6	0.57	0
2	SO4	A	463	-	4,4,4	1.05	0	6,6,6	0.58	0
2	SO4	C	436	-	4,4,4	1.06	0	6,6,6	0.58	0
2	SO4	B	449	-	4,4,4	1.07	0	6,6,6	0.58	0
2	SO4	D	442	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	C	444	-	4,4,4	1.07	0	6,6,6	0.58	0
2	SO4	B	434	-	4,4,4	1.10	0	6,6,6	0.56	0
2	SO4	A	456	-	4,4,4	1.09	0	6,6,6	0.57	0
2	SO4	A	457	-	4,4,4	1.05	0	6,6,6	0.58	0
2	SO4	A	459	-	4,4,4	1.05	0	6,6,6	0.58	0
2	SO4	C	456	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	A	444	-	4,4,4	1.03	0	6,6,6	0.59	0
2	SO4	B	444	-	4,4,4	1.07	0	6,6,6	0.56	0
2	SO4	C	448	-	4,4,4	1.07	0	6,6,6	0.56	0
2	SO4	D	447	-	4,4,4	1.07	0	6,6,6	0.58	0
2	SO4	A	443	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	C	435	-	4,4,4	1.06	0	6,6,6	0.56	0
2	SO4	D	445	-	4,4,4	1.07	0	6,6,6	0.57	0
2	SO4	D	433	-	4,4,4	1.06	0	6,6,6	0.57	0
2	SO4	B	432	-	4,4,4	1.06	0	6,6,6	0.56	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	U5P	D	448	-	-	0/10/26/26	0/2/2/2
4	U5P	A	467	-	-	3/10/26/26	0/2/2/2
4	U5P	C	457	-	-	0/10/26/26	0/2/2/2
4	U5P	B	457	-	-	6/10/26/26	0/2/2/2
4	U5P	B	456	-	-	2/10/26/26	0/2/2/2
3	FLC	A	464	-	-	0/16/16/16	-
4	U5P	A	465	-	-	5/10/26/26	0/2/2/2
4	U5P	A	466	-	-	4/10/26/26	0/2/2/2
3	FLC	B	455	-	-	0/16/16/16	-

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	457	U5P	C2-N1	6.80	1.49	1.38
4	A	466	U5P	C2-N1	6.08	1.48	1.38
4	A	467	U5P	C2-N1	6.07	1.48	1.38
4	A	465	U5P	C2-N1	6.01	1.47	1.38
4	D	448	U5P	C2-N1	6.01	1.47	1.38
4	B	456	U5P	C2-N1	6.00	1.47	1.38
4	C	457	U5P	C2-N1	5.57	1.47	1.38
4	A	466	U5P	C6-C5	4.03	1.44	1.35
4	B	457	U5P	C6-C5	3.96	1.44	1.35
4	A	467	U5P	C6-C5	3.93	1.44	1.35
4	C	457	U5P	C6-C5	3.92	1.44	1.35
4	D	448	U5P	C6-C5	3.85	1.44	1.35
4	B	456	U5P	C6-C5	3.71	1.43	1.35
4	A	465	U5P	C6-C5	3.63	1.43	1.35
4	A	466	U5P	O2-C2	3.31	1.28	1.23
4	A	467	U5P	O2-C2	3.25	1.28	1.23
4	C	457	U5P	P-O5'	-3.17	1.50	1.60
4	D	448	U5P	O2-C2	3.16	1.28	1.23
4	C	457	U5P	O2-C2	3.13	1.28	1.23
4	A	465	U5P	P-O5'	-3.06	1.50	1.60
4	D	448	U5P	P-O5'	-3.04	1.50	1.60
4	B	456	U5P	P-O5'	-2.97	1.50	1.60
4	A	465	U5P	O2-C2	2.96	1.28	1.23
4	B	457	U5P	O2-C2	2.94	1.28	1.23
4	B	456	U5P	O2-C2	2.93	1.28	1.23
4	A	467	U5P	P-O5'	-2.89	1.51	1.60
4	B	457	U5P	P-O5'	-2.72	1.51	1.60
3	B	455	FLC	CG-CB	2.69	1.57	1.54
4	A	466	U5P	P-O5'	-2.60	1.52	1.60
3	A	464	FLC	OG2-CGC	-2.49	1.22	1.30
3	A	464	FLC	OA2-CAC	-2.46	1.22	1.30
4	A	466	U5P	P-O2P	2.44	1.63	1.54
3	B	455	FLC	OA2-CAC	-2.42	1.22	1.30
4	C	457	U5P	P-O2P	2.36	1.63	1.54
3	B	455	FLC	OG2-CGC	-2.34	1.23	1.30
4	B	457	U5P	P-O2P	2.34	1.63	1.54
3	B	455	FLC	CA-CB	2.32	1.56	1.54
4	D	448	U5P	P-O2P	2.24	1.63	1.54
4	A	467	U5P	P-O2P	2.20	1.63	1.54
3	A	464	FLC	CA-CB	2.14	1.56	1.54
3	B	455	FLC	OB1-CBC	2.10	1.28	1.22
3	A	464	FLC	OB1-CBC	2.10	1.28	1.22
3	A	464	FLC	CG-CB	2.03	1.56	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	465	U5P	P-O2P	2.02	1.62	1.54

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	448	U5P	C5-C4-N3	4.98	121.77	114.80
4	C	457	U5P	C5-C4-N3	4.95	121.74	114.80
4	A	465	U5P	C5-C4-N3	4.87	121.62	114.80
4	A	467	U5P	C5-C4-N3	4.87	121.62	114.80
4	B	456	U5P	C5-C4-N3	4.85	121.60	114.80
4	B	457	U5P	C5-C4-N3	4.78	121.50	114.80
4	A	466	U5P	C5-C4-N3	4.66	121.33	114.80
4	A	465	U5P	C4-N3-C2	-4.13	121.49	126.61
4	B	456	U5P	C4-N3-C2	-4.01	121.63	126.61
3	A	464	FLC	OB2-CBC-CB	3.83	120.50	113.14
4	C	457	U5P	C4-N3-C2	-3.79	121.90	126.61
4	D	448	U5P	C4-N3-C2	-3.75	121.96	126.61
4	A	467	U5P	C4-N3-C2	-3.70	122.02	126.61
4	B	457	U5P	C4-N3-C2	-3.55	122.21	126.61
4	A	466	U5P	C4-N3-C2	-3.53	122.23	126.61
3	B	455	FLC	OB2-CBC-CB	3.51	119.88	113.14
4	A	467	U5P	C2'-C1'-N1	3.32	122.48	113.25
4	A	465	U5P	N3-C2-N1	3.06	118.88	114.89
4	A	465	U5P	C2'-C1'-N1	3.06	121.77	113.25
4	D	448	U5P	C2'-C1'-N1	3.05	121.74	113.25
4	B	457	U5P	N3-C2-N1	3.02	118.83	114.89
4	A	466	U5P	C2'-C1'-N1	2.96	121.48	113.25
4	B	456	U5P	N3-C2-N1	2.94	118.72	114.89
4	C	457	U5P	N3-C2-N1	2.94	118.71	114.89
4	B	456	U5P	C2'-C1'-N1	2.92	121.38	113.25
4	A	466	U5P	N3-C2-N1	2.89	118.66	114.89
4	A	467	U5P	N3-C2-N1	2.89	118.65	114.89
4	D	448	U5P	N3-C2-N1	2.88	118.64	114.89
4	B	457	U5P	C2'-C1'-N1	2.76	120.93	113.25
4	A	465	U5P	C1'-N1-C6	2.73	126.61	120.78
4	C	457	U5P	C2'-C1'-N1	2.70	120.77	113.25
4	A	467	U5P	C1'-N1-C6	2.61	126.36	120.78
4	C	457	U5P	C1'-N1-C6	2.55	126.22	120.78
4	A	466	U5P	C1'-N1-C6	2.46	126.03	120.78
4	D	448	U5P	C1'-N1-C6	2.43	125.98	120.78
4	B	456	U5P	C1'-N1-C6	2.38	125.87	120.78
4	A	465	U5P	O4-C4-C5	-2.28	121.23	125.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	456	U5P	O4-C4-C5	-2.23	121.31	125.16
4	B	457	U5P	C6-N1-C2	-2.15	118.38	121.00
4	B	457	U5P	O2-C2-N3	-2.05	117.70	121.49
4	A	466	U5P	O4-C4-C5	-2.04	121.64	125.16
4	C	457	U5P	O4-C4-C5	-2.01	121.69	125.16

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	465	U5P	C5'-O5'-P-O2P
4	A	465	U5P	C5'-O5'-P-O3P
4	A	466	U5P	C5'-O5'-P-O2P
4	A	466	U5P	C5'-O5'-P-O3P
4	B	457	U5P	O4'-C1'-N1-C2
4	B	457	U5P	O4'-C1'-N1-C6
4	B	457	U5P	C5'-O5'-P-O2P
4	B	457	U5P	C5'-O5'-P-O3P
4	B	457	U5P	C3'-C4'-C5'-O5'
4	B	457	U5P	O4'-C4'-C5'-O5'
4	A	465	U5P	O4'-C4'-C5'-O5'
4	A	465	U5P	C5'-O5'-P-O1P
4	A	465	U5P	C3'-C4'-C5'-O5'
4	A	467	U5P	C5'-O5'-P-O2P
4	A	466	U5P	C3'-C4'-C5'-O5'
4	A	466	U5P	O4'-C4'-C5'-O5'
4	A	467	U5P	C5'-O5'-P-O1P
4	B	456	U5P	C5'-O5'-P-O1P
4	B	456	U5P	C5'-O5'-P-O2P
4	A	467	U5P	C4'-C5'-O5'-P

There are no ring outliers.

16 monomers are involved in 22 short contacts:

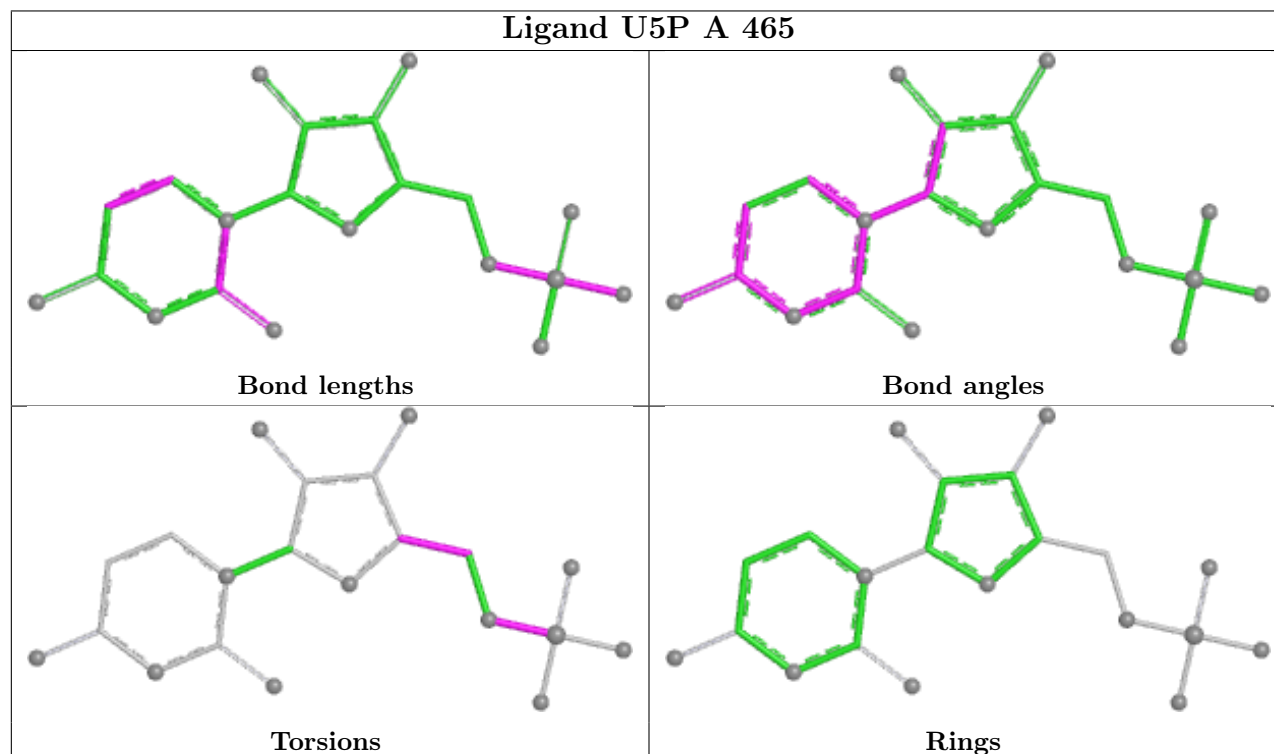
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	465	U5P	2	0
2	D	443	SO4	1	0
2	A	433	SO4	3	0
3	B	455	FLC	2	0
4	B	456	U5P	1	0
4	B	457	U5P	1	0

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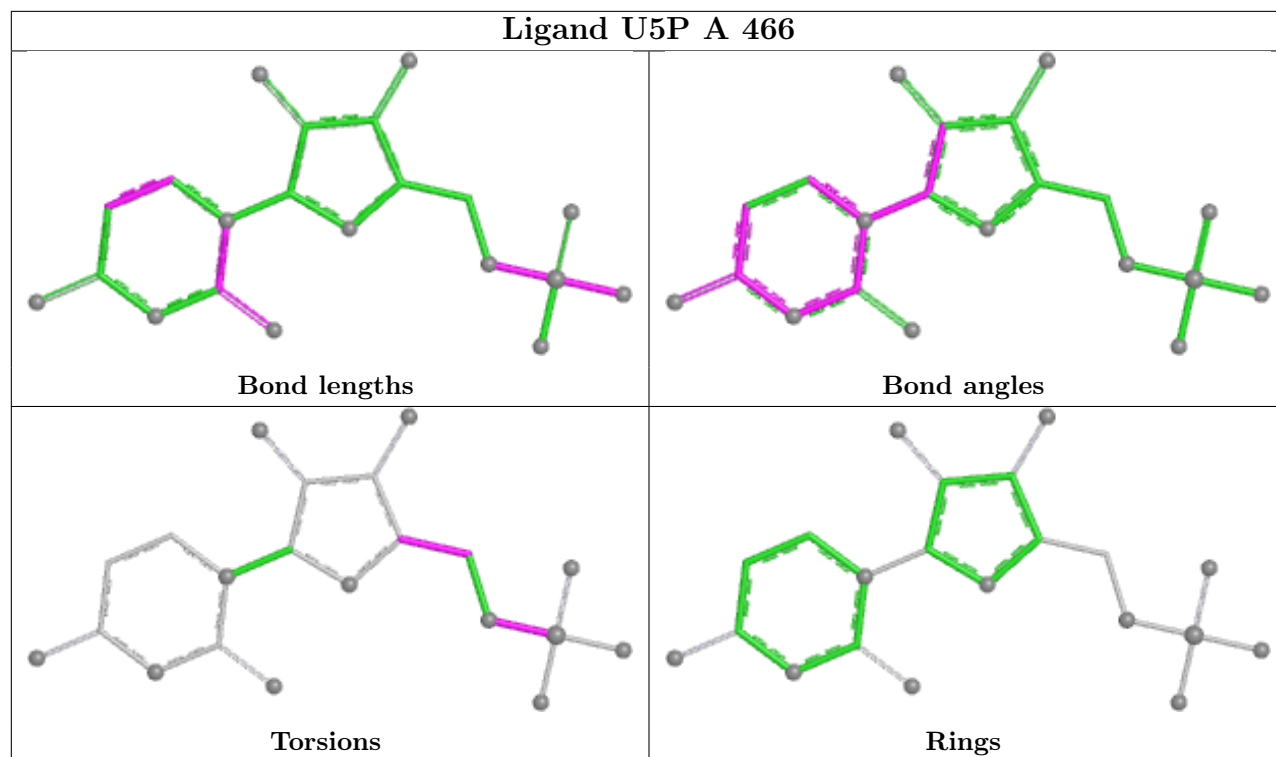
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	464	FLC	1	0
2	C	453	SO4	2	0
2	A	451	SO4	1	0
2	C	451	SO4	1	0
2	C	433	SO4	1	0
2	C	454	SO4	1	0
2	A	463	SO4	1	0
2	A	459	SO4	1	0
2	A	444	SO4	1	0
2	A	443	SO4	2	0

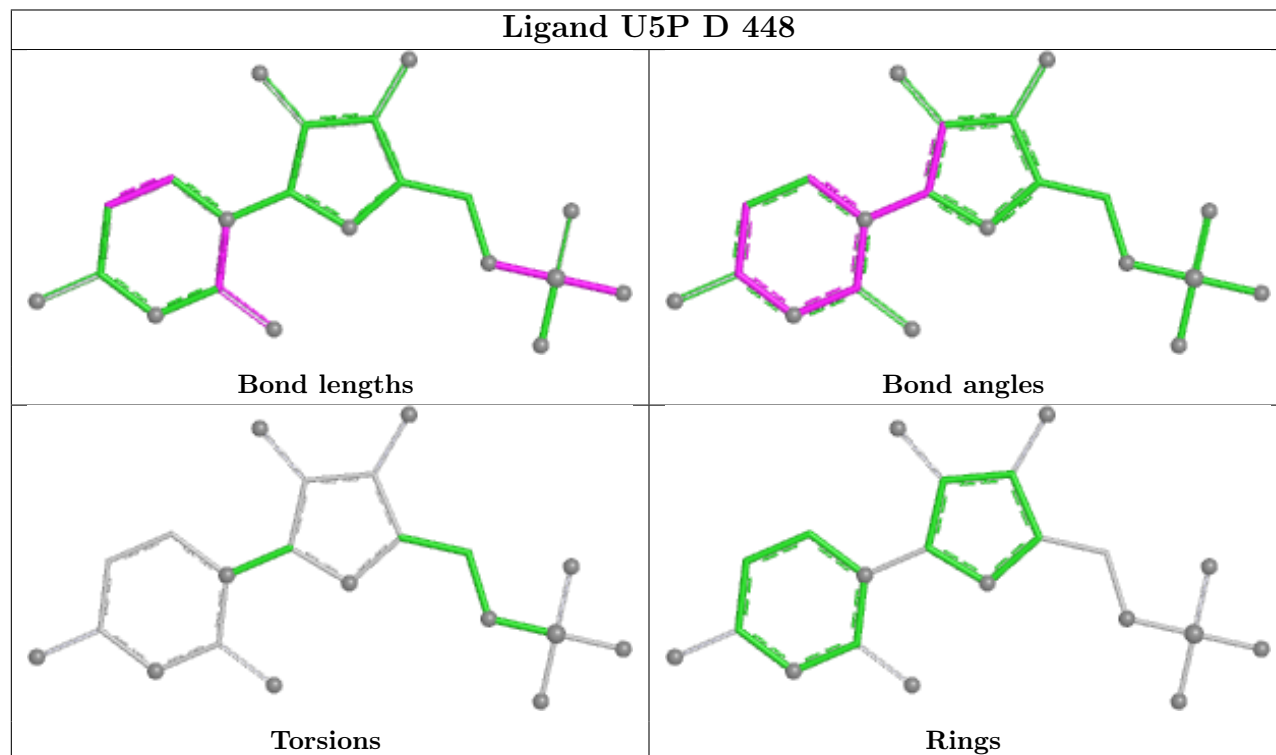
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.



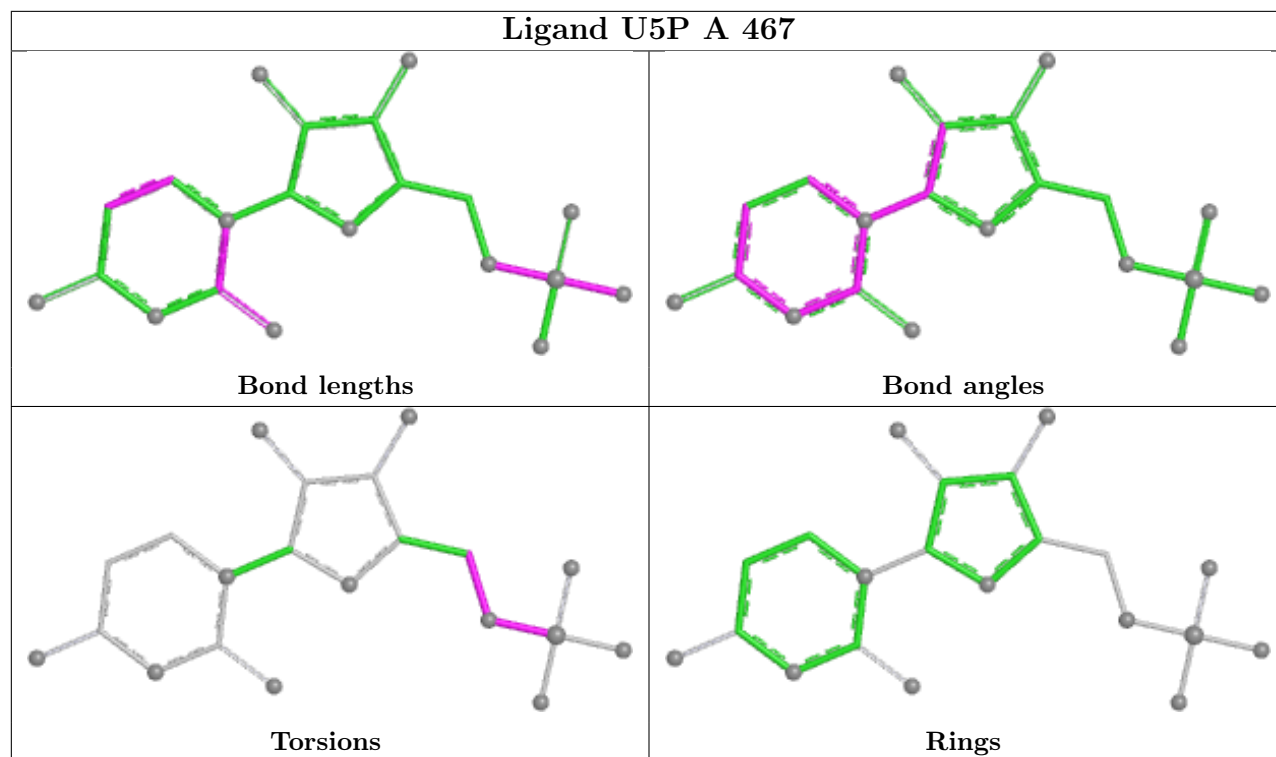
Ligand U5P A 466



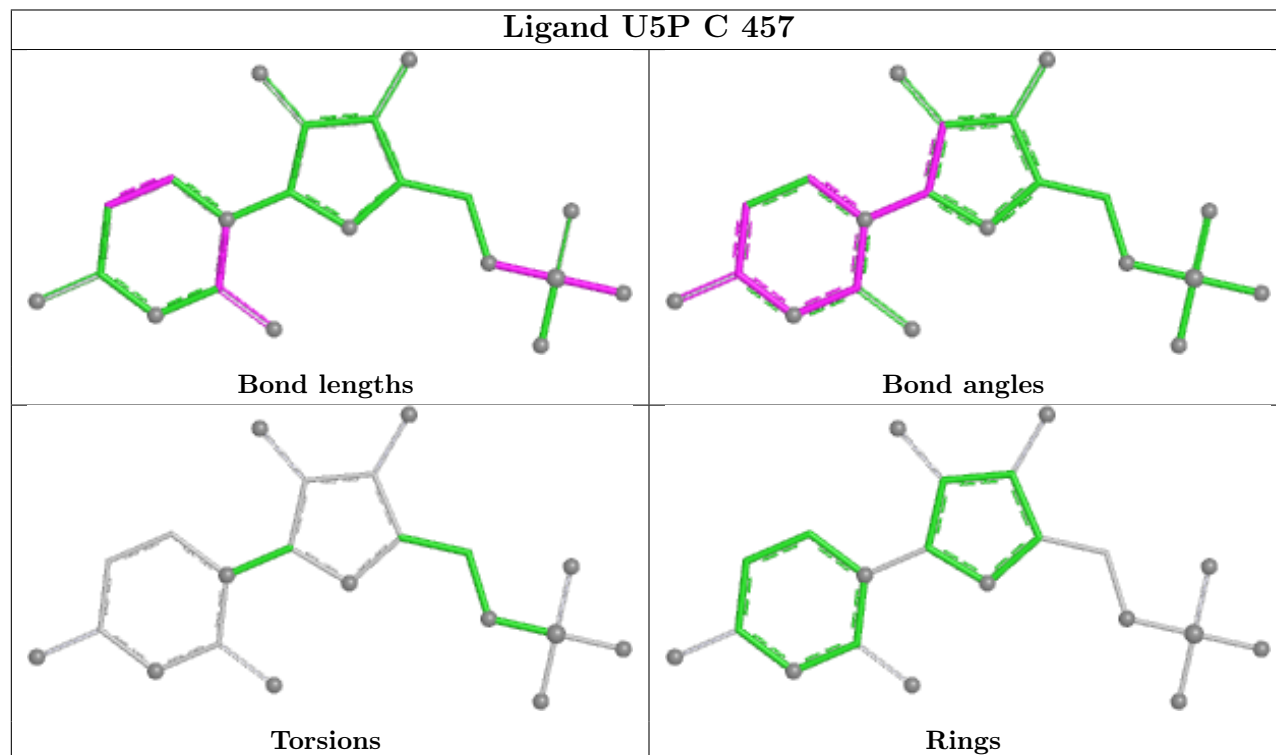
Ligand U5P D 448

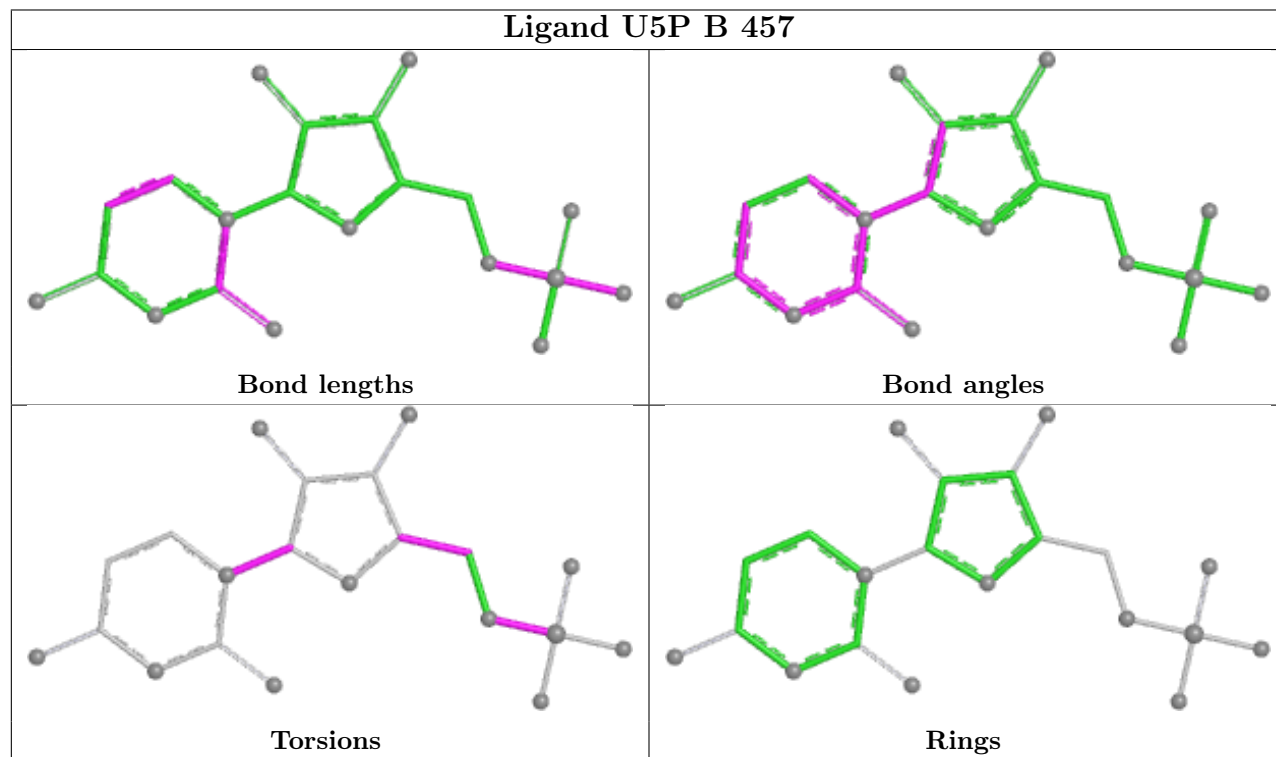
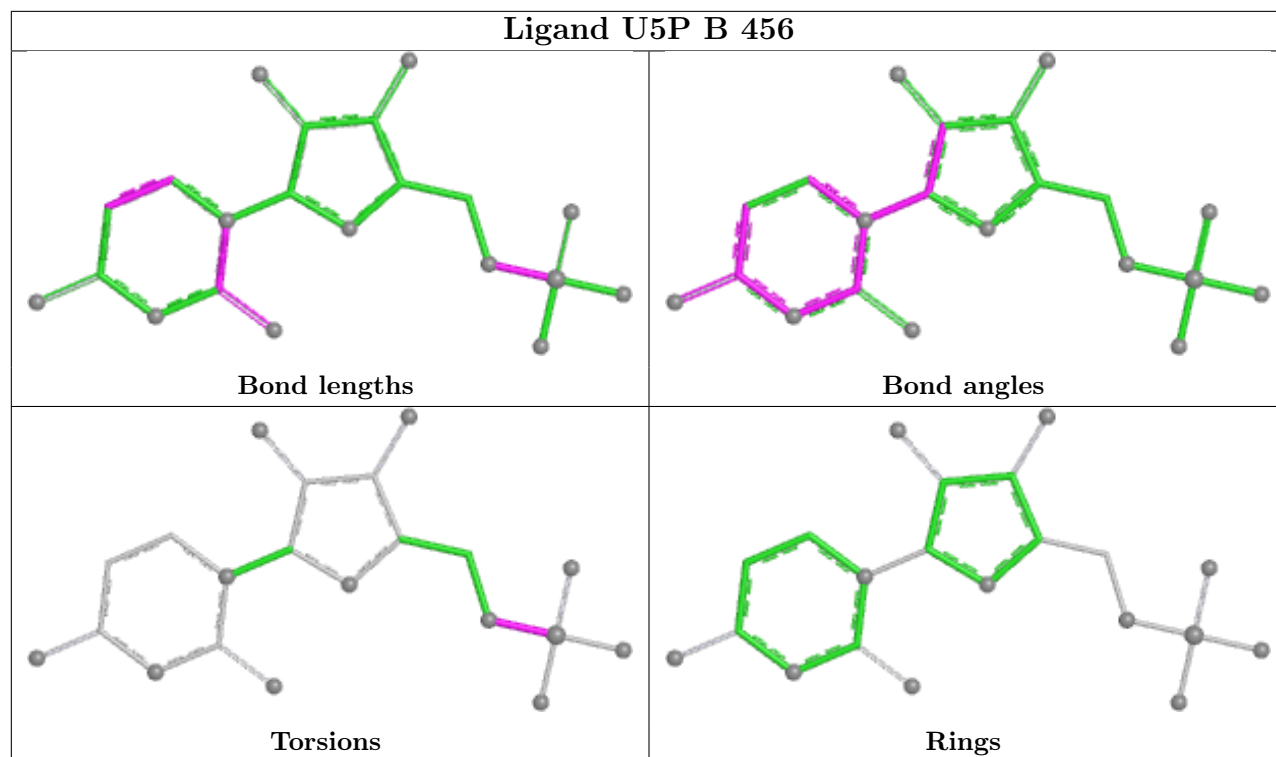


Ligand U5P A 467



Ligand U5P C 457





5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	431/431 (100%)	0.08	7 (1%) 70 76	26, 39, 62, 83	0
1	B	431/431 (100%)	0.21	11 (2%) 57 63	23, 40, 64, 83	0
1	C	431/431 (100%)	1.01	81 (18%) 3 3	26, 61, 109, 121	0
1	D	431/431 (100%)	1.48	124 (28%) 1 1	36, 69, 124, 144	0
All	All	1724/1724 (100%)	0.70	223 (12%) 7 9	23, 48, 111, 144	0

All (223) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	365	VAL	6.1
1	D	339	VAL	5.5
1	D	358	VAL	5.5
1	D	362	GLY	5.1
1	D	300	LEU	4.9
1	D	366	PRO	4.8
1	D	299	ALA	4.8
1	D	376	GLY	4.8
1	D	327	HIS	4.6
1	D	340	GLY	4.5
1	D	325	LEU	4.4
1	D	249	LEU	4.4
1	D	242	LEU	4.4
1	D	338	PHE	4.3
1	D	219	LEU	4.2
1	C	303	ALA	4.2
1	D	363	GLU	4.1
1	D	341	TYR	4.0
1	D	357	ALA	4.0
1	D	351	ILE	4.0
1	D	361	LEU	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	291	VAL	3.9
1	D	329	LEU	3.9
1	D	328	GLY	3.9
1	D	212	LEU	3.8
1	C	335	ALA	3.8
1	D	369	ALA	3.8
1	D	324	HIS	3.8
1	C	277	LEU	3.8
1	C	291	VAL	3.7
1	D	208	LEU	3.7
1	D	334	ASN	3.7
1	D	367	LEU	3.7
1	D	360	ILE	3.6
1	D	250	ASP	3.6
1	C	246	PRO	3.6
1	D	201	VAL	3.6
1	D	314	GLY	3.6
1	D	306	PRO	3.6
1	D	200	THR	3.6
1	D	332	PRO	3.5
1	D	253	MET	3.5
1	D	309	VAL	3.4
1	D	296	ALA	3.4
1	D	252	PRO	3.4
1	D	100	VAL	3.4
1	C	216	GLY	3.4
1	D	355	PRO	3.4
1	C	370	SER	3.4
1	C	359	ARG	3.4
1	C	288	LEU	3.4
1	D	205	LEU	3.4
1	D	225	VAL	3.4
1	C	301	ASN	3.4
1	D	415	LEU	3.3
1	D	204	PHE	3.3
1	D	302	ARG	3.3
1	D	293	HIS	3.3
1	C	274	ALA	3.3
1	D	311	ALA	3.3
1	C	308	VAL	3.3
1	D	335	ALA	3.3
1	D	356	PRO	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	39	ALA	3.2
1	C	236	TYR	3.2
1	C	265	VAL	3.2
1	D	248	TYR	3.2
1	C	333	ARG	3.1
1	D	286	ALA	3.1
1	D	223	PHE	3.1
1	D	377	PHE	3.1
1	D	101	MET	3.1
1	D	303	ALA	3.1
1	D	336	LEU	3.1
1	C	307	MET	3.1
1	D	347	LEU	3.0
1	A	302	ARG	3.0
1	D	298	LYS	3.0
1	D	310	LEU	3.0
1	D	247	ILE	3.0
1	D	331	ASP	3.0
1	D	317	ALA	3.0
1	C	100	VAL	3.0
1	D	371	VAL	3.0
1	D	297	SER	3.0
1	D	102	ASP	3.0
1	C	286	ALA	2.9
1	D	104	PRO	2.9
1	D	207	ILE	2.9
1	B	23	GLY	2.9
1	C	357	ALA	2.9
1	D	98	LEU	2.9
1	D	431	VAL	2.9
1	B	22	GLY	2.8
1	D	375	GLY	2.8
1	C	299	ALA	2.8
1	D	337	VAL	2.8
1	C	352	ILE	2.8
1	C	367	LEU	2.8
1	C	276	PHE	2.8
1	D	307	MET	2.8
1	B	240	HIS	2.8
1	D	305	GLY	2.8
1	D	285	PRO	2.8
1	C	306	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	37	GLU	2.7
1	C	242	LEU	2.7
1	C	358	VAL	2.7
1	D	326	LYS	2.7
1	C	300	LEU	2.7
1	D	235	LEU	2.7
1	C	283	PHE	2.7
1	D	210	LYS	2.6
1	C	365	VAL	2.6
1	D	301	ASN	2.6
1	C	305	GLY	2.6
1	D	240	HIS	2.6
1	C	296	ALA	2.6
1	D	313	SER	2.6
1	C	336	LEU	2.6
1	D	352	ILE	2.6
1	D	372	HIS	2.6
1	C	202	ARG	2.6
1	B	43	ALA	2.6
1	D	243	PRO	2.6
1	D	231	ILE	2.6
1	C	329	LEU	2.5
1	C	302	ARG	2.5
1	D	244	ARG	2.5
1	D	320	ARG	2.5
1	C	311	ALA	2.5
1	D	211	THR	2.5
1	D	222	THR	2.5
1	C	225	VAL	2.5
1	C	233	TYR	2.5
1	B	38	GLU	2.5
1	C	219	LEU	2.5
1	C	353	ALA	2.5
1	C	215	GLY	2.5
1	C	297	SER	2.5
1	C	330	SER	2.5
1	C	247	ILE	2.5
1	C	369	ALA	2.5
1	C	221	PRO	2.5
1	C	279	GLY	2.5
1	C	356	PRO	2.5
1	D	188	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	308	VAL	2.5
1	C	304	PRO	2.5
1	C	364	GLU	2.4
1	D	105	PHE	2.4
1	D	374	LEU	2.4
1	D	239	GLY	2.4
1	C	275	HIS	2.4
1	B	103	GLU	2.4
1	B	404	GLU	2.4
1	D	246	PRO	2.4
1	D	318	GLY	2.4
1	D	213	SER	2.4
1	C	285	PRO	2.4
1	C	245	ALA	2.4
1	C	237	THR	2.4
1	D	373	THR	2.4
1	D	218	VAL	2.4
1	D	220	ILE	2.3
1	C	266	ARG	2.3
1	D	292	GLU	2.3
1	C	240	HIS	2.3
1	D	364	GLU	2.3
1	D	316	LEU	2.3
1	D	197	TYR	2.3
1	A	365	VAL	2.3
1	A	343	PRO	2.3
1	C	243	PRO	2.3
1	D	354	ARG	2.3
1	D	217	LYS	2.3
1	A	129	LEU	2.2
1	C	208	LEU	2.2
1	C	223	PHE	2.2
1	C	332	PRO	2.2
1	C	293	HIS	2.2
1	D	330	SER	2.2
1	D	344	GLN	2.2
1	C	207	ILE	2.2
1	C	351	ILE	2.2
1	D	321	ILE	2.2
1	A	100	VAL	2.2
1	D	167	GLU	2.2
1	C	258	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	241	ARG	2.2
1	D	94	LEU	2.2
1	C	371	VAL	2.2
1	C	295	GLU	2.2
1	C	290	VAL	2.1
1	D	290	VAL	2.1
1	D	370	SER	2.1
1	C	171	LEU	2.1
1	C	292	GLU	2.1
1	D	283	PHE	2.1
1	C	375	GLY	2.1
1	D	312	GLY	2.1
1	D	368	ARG	2.1
1	D	353	ALA	2.1
1	C	280	LYS	2.1
1	C	326	LYS	2.1
1	C	361	LEU	2.1
1	D	236	TYR	2.1
1	A	225	VAL	2.1
1	D	359	ARG	2.1
1	B	42	HIS	2.1
1	D	214	GLN	2.1
1	C	339	VAL	2.1
1	D	170	VAL	2.1
1	D	251	SER	2.0
1	B	101	MET	2.0
1	C	98	LEU	2.0
1	C	412	LEU	2.0
1	D	281	ASN	2.0
1	C	201	VAL	2.0
1	A	362	GLY	2.0
1	C	23	GLY	2.0
1	C	373	THR	2.0
1	C	324	HIS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	C	455	5/5	0.35	0.14	138,138,139,139	0
2	SO4	D	438	5/5	0.35	0.12	159,159,159,159	0
2	SO4	D	443	5/5	0.44	0.15	185,185,185,185	0
2	SO4	C	452	5/5	0.46	0.11	149,149,149,149	0
2	SO4	C	442	5/5	0.46	0.15	150,151,151,151	0
2	SO4	D	441	5/5	0.47	0.21	173,173,173,174	0
2	SO4	A	436	5/5	0.50	0.15	159,159,159,159	0
2	SO4	C	433	5/5	0.50	0.17	138,138,138,138	0
2	SO4	A	450	5/5	0.52	0.15	166,166,166,166	0
2	SO4	A	461	5/5	0.53	0.14	125,125,126,126	0
2	SO4	B	438	5/5	0.54	0.18	133,133,133,133	0
2	SO4	A	447	5/5	0.54	0.20	145,145,145,146	0
2	SO4	A	433	5/5	0.55	0.21	157,158,158,158	0
2	SO4	D	446	5/5	0.55	0.12	165,165,166,166	0
2	SO4	D	445	5/5	0.57	0.15	131,131,132,132	0
2	SO4	D	447	5/5	0.57	0.13	161,161,161,161	0
2	SO4	C	446	5/5	0.58	0.11	155,155,155,156	0
2	SO4	A	442	5/5	0.59	0.19	159,159,159,159	0
2	SO4	C	453	5/5	0.59	0.18	152,152,152,152	0
2	SO4	C	449	5/5	0.59	0.09	148,148,148,148	0
4	U5P	A	467	21/21	0.60	0.19	140,144,145,146	0
2	SO4	C	447	5/5	0.61	0.11	186,186,186,186	0
2	SO4	B	450	5/5	0.61	0.18	122,122,123,123	0
2	SO4	C	444	5/5	0.62	0.12	124,124,124,124	0
2	SO4	A	458	5/5	0.62	0.10	124,124,125,125	0
2	SO4	D	433	5/5	0.63	0.18	141,141,142,142	0
2	SO4	C	441	5/5	0.63	0.10	130,130,130,130	0
2	SO4	B	451	5/5	0.63	0.13	143,143,143,143	0
2	SO4	A	444	5/5	0.63	0.13	138,138,138,138	0
2	SO4	C	456	5/5	0.64	0.13	136,137,137,137	0
2	SO4	D	442	5/5	0.65	0.21	167,167,167,167	0
2	SO4	A	445	5/5	0.65	0.15	125,125,125,126	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	A	434	5/5	0.66	0.18	140,140,140,140	0
2	SO4	B	435	5/5	0.66	0.14	132,132,133,133	0
2	SO4	A	449	5/5	0.66	0.12	156,156,156,156	0
2	SO4	A	432	5/5	0.66	0.15	123,123,124,124	0
2	SO4	A	439	5/5	0.66	0.17	131,131,131,131	0
2	SO4	A	443	5/5	0.67	0.16	141,141,142,142	0
2	SO4	D	434	5/5	0.67	0.12	123,123,123,123	0
2	SO4	C	454	5/5	0.67	0.11	143,144,144,144	0
2	SO4	C	436	5/5	0.67	0.12	129,130,130,130	0
2	SO4	A	451	5/5	0.67	0.18	130,130,130,130	0
2	SO4	A	463	5/5	0.68	0.16	142,142,143,143	0
2	SO4	C	438	5/5	0.69	0.13	123,123,124,124	0
2	SO4	A	446	5/5	0.69	0.21	167,167,167,167	0
4	U5P	B	457	21/21	0.69	0.19	106,108,110,111	0
2	SO4	D	437	5/5	0.70	0.15	141,141,142,142	0
2	SO4	D	440	5/5	0.70	0.09	139,139,139,139	0
2	SO4	A	435	5/5	0.71	0.16	126,126,127,127	0
2	SO4	A	438	5/5	0.71	0.15	143,144,144,144	0
2	SO4	B	448	5/5	0.72	0.12	105,105,106,106	0
3	FLC	B	455	13/13	0.72	0.23	84,97,99,100	0
2	SO4	A	448	5/5	0.72	0.19	125,125,125,125	0
2	SO4	A	460	5/5	0.72	0.15	101,101,101,102	0
2	SO4	B	439	5/5	0.73	0.14	135,135,135,135	0
2	SO4	D	439	5/5	0.73	0.11	131,131,132,132	0
2	SO4	C	432	5/5	0.73	0.14	108,108,108,108	0
2	SO4	C	437	5/5	0.73	0.09	142,142,142,142	0
2	SO4	B	440	5/5	0.74	0.21	140,140,140,141	0
2	SO4	B	447	5/5	0.74	0.15	142,142,142,142	0
2	SO4	B	437	5/5	0.75	0.11	119,119,119,119	0
2	SO4	A	459	5/5	0.75	0.15	118,119,119,119	0
2	SO4	C	448	5/5	0.75	0.14	141,141,141,141	0
2	SO4	B	441	5/5	0.76	0.12	107,107,108,109	0
2	SO4	A	452	5/5	0.76	0.13	120,120,120,120	0
2	SO4	C	440	5/5	0.76	0.21	174,174,174,174	0
2	SO4	A	457	5/5	0.76	0.18	129,129,129,130	0
2	SO4	A	462	5/5	0.76	0.15	127,128,128,128	0
2	SO4	D	432	5/5	0.77	0.14	132,132,132,132	0
2	SO4	B	454	5/5	0.77	0.17	147,147,148,148	0
2	SO4	D	436	5/5	0.78	0.09	108,108,108,108	0
2	SO4	C	451	5/5	0.78	0.12	111,111,111,111	0
2	SO4	C	439	5/5	0.79	0.12	111,111,112,112	0
2	SO4	B	452	5/5	0.79	0.15	103,103,103,104	0

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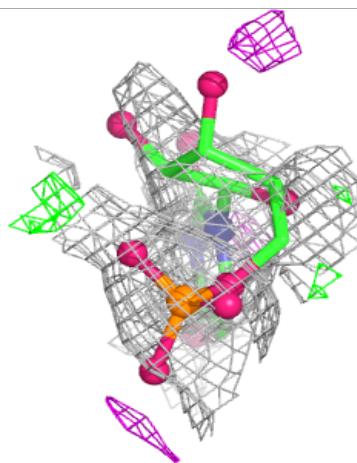
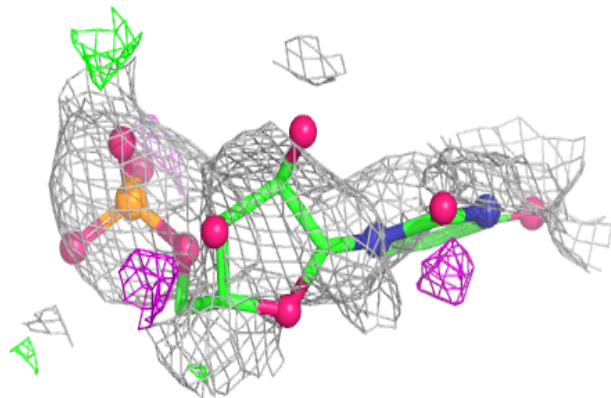
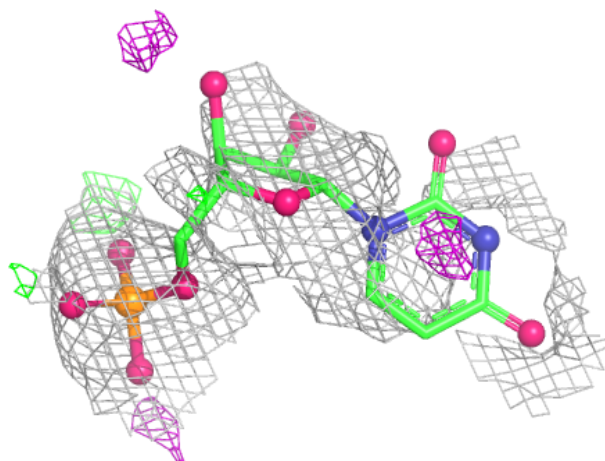
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	C	450	5/5	0.80	0.13	142,142,143,143	0
2	SO4	B	453	5/5	0.80	0.13	124,124,124,125	0
2	SO4	C	435	5/5	0.81	0.12	113,114,114,114	0
2	SO4	B	442	5/5	0.81	0.14	106,106,107,107	0
2	SO4	C	443	5/5	0.81	0.13	117,117,118,118	0
2	SO4	A	456	5/5	0.81	0.13	90,90,91,91	0
3	FLC	A	464	13/13	0.82	0.26	125,127,128,128	0
4	U5P	A	466	21/21	0.82	0.14	60,71,74,76	0
2	SO4	B	449	5/5	0.83	0.15	141,141,141,141	0
2	SO4	A	441	5/5	0.85	0.10	101,101,101,101	0
2	SO4	A	455	5/5	0.85	0.14	114,114,114,114	0
2	SO4	B	434	5/5	0.85	0.14	84,84,85,85	0
2	SO4	B	444	5/5	0.86	0.12	109,110,110,110	0
2	SO4	C	445	5/5	0.87	0.14	99,99,100,100	0
4	U5P	D	448	21/21	0.87	0.14	88,94,95,96	0
2	SO4	A	437	5/5	0.88	0.11	87,87,88,89	0
5	ZN	C	458	1/1	0.88	0.16	96,96,96,96	0
2	SO4	B	446	5/5	0.89	0.14	135,135,135,136	0
2	SO4	B	443	5/5	0.90	0.09	80,81,82,82	0
2	SO4	C	434	5/5	0.90	0.14	70,71,71,72	0
2	SO4	B	436	5/5	0.90	0.21	94,95,95,97	0
2	SO4	A	453	5/5	0.91	0.09	84,85,85,86	0
2	SO4	D	444	5/5	0.91	0.13	72,72,73,73	0
2	SO4	D	435	5/5	0.91	0.11	87,87,88,88	0
2	SO4	A	454	5/5	0.91	0.12	54,54,55,57	0
4	U5P	C	457	21/21	0.92	0.12	61,73,75,76	0
2	SO4	B	432	5/5	0.93	0.09	83,84,85,86	0
5	ZN	C	459	1/1	0.93	0.10	97,97,97,97	0
2	SO4	B	445	5/5	0.94	0.11	58,60,62,63	0
5	ZN	A	469	1/1	0.94	0.12	95,95,95,95	0
5	ZN	D	449	1/1	0.94	0.12	106,106,106,106	0
5	ZN	B	458	1/1	0.95	0.14	70,70,70,70	0
4	U5P	A	465	21/21	0.95	0.10	33,51,52,57	0
4	U5P	B	456	21/21	0.95	0.10	34,53,56,59	0
2	SO4	B	433	5/5	0.95	0.08	65,65,66,66	0
5	ZN	B	459	1/1	0.96	0.13	84,84,84,84	0
2	SO4	A	440	5/5	0.96	0.08	58,58,60,60	0
5	ZN	D	450	1/1	0.96	0.13	102,102,102,102	0
5	ZN	A	468	1/1	0.97	0.15	74,74,74,74	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.

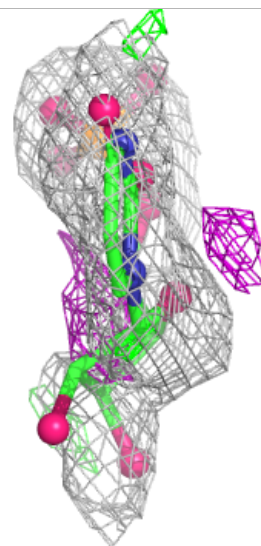
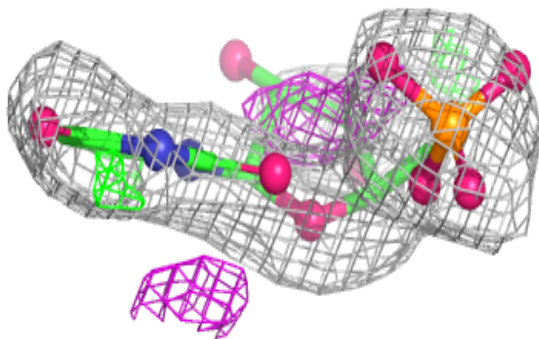
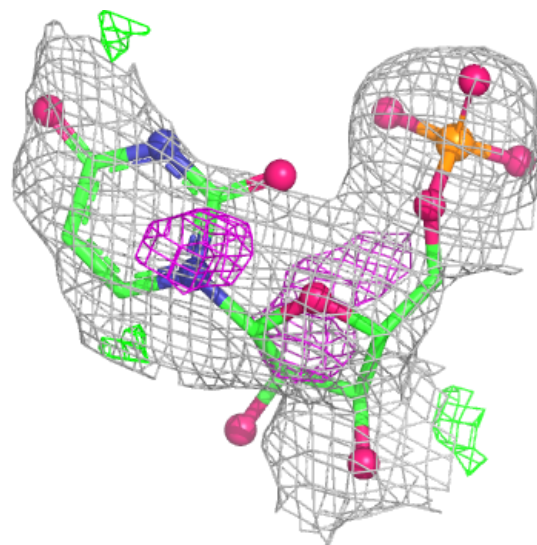
Electron density around U5P A 467:

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



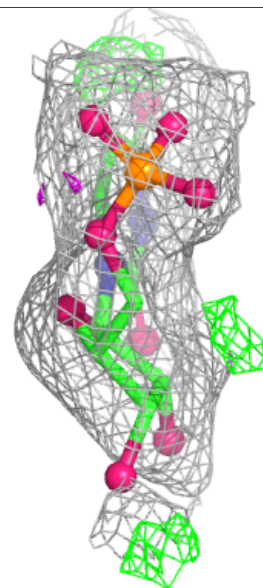
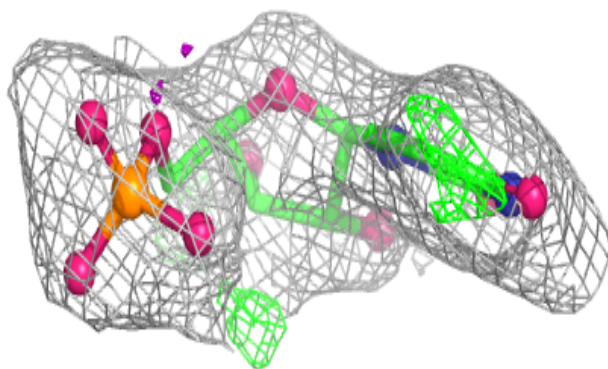
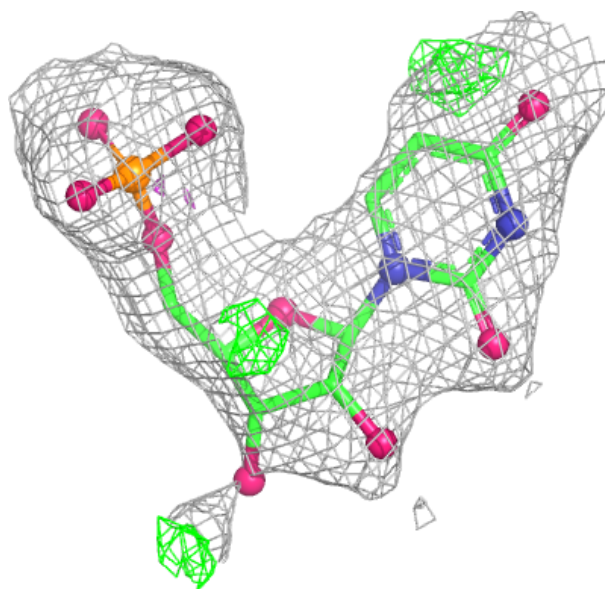
Electron density around U5P B 457:

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



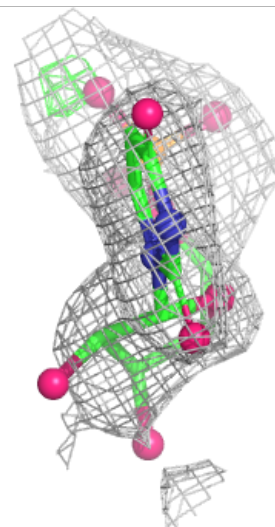
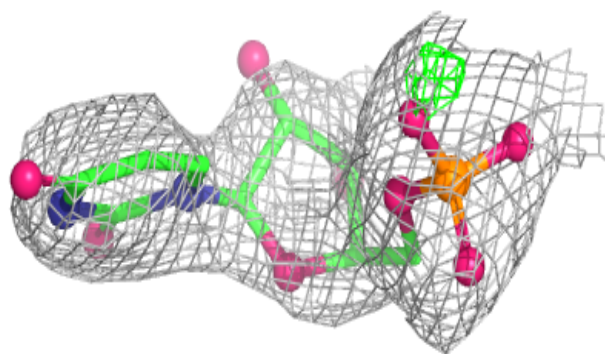
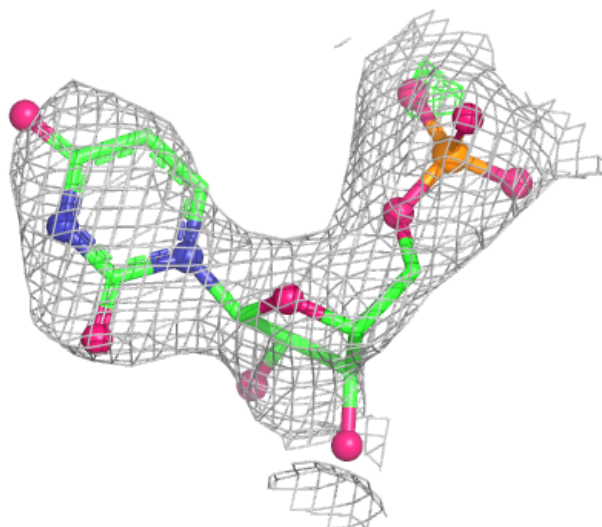
Electron density around U5P A 466:

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



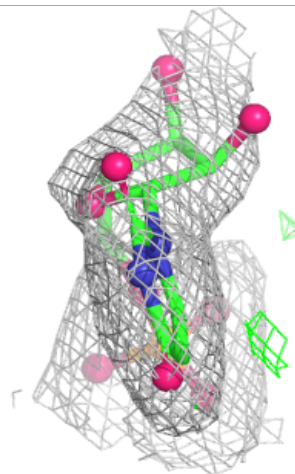
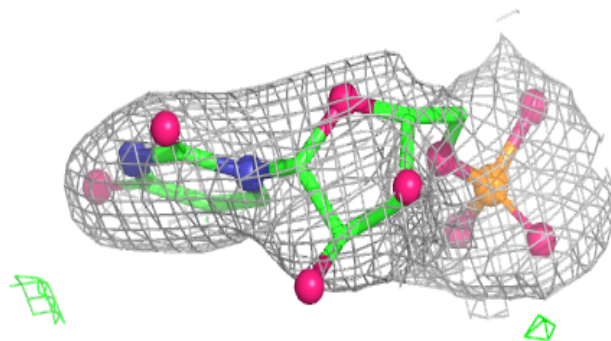
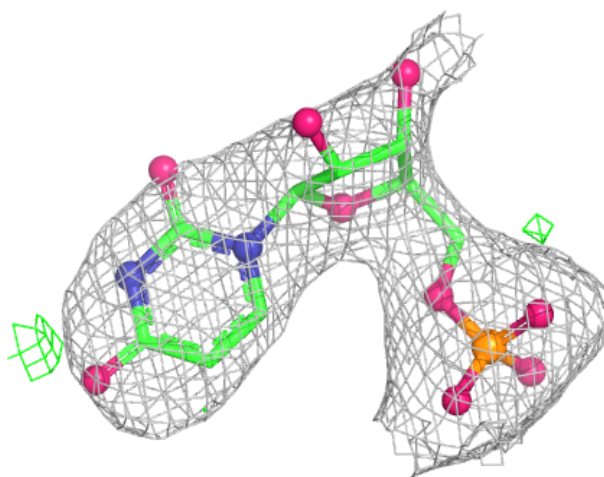
Electron density around U5P D 448:

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



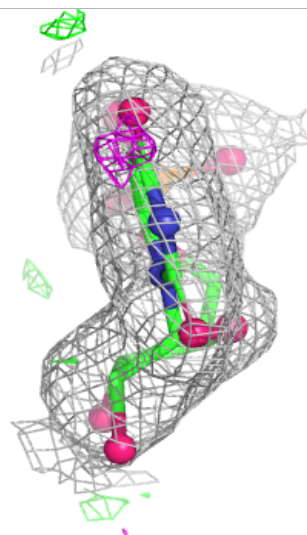
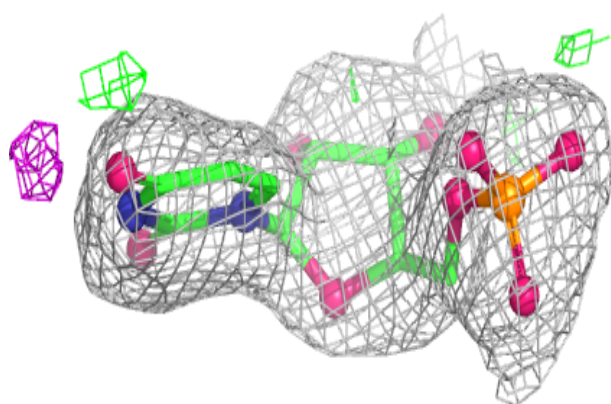
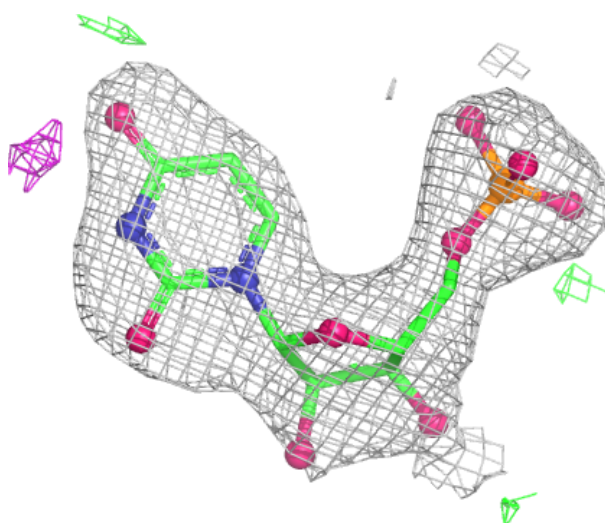
Electron density around U5P C 457:

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



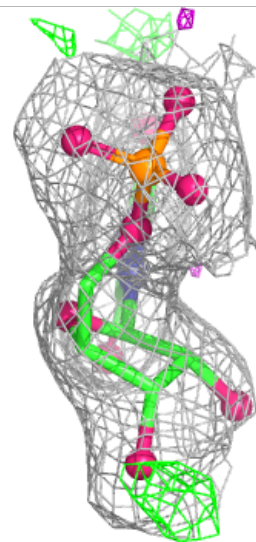
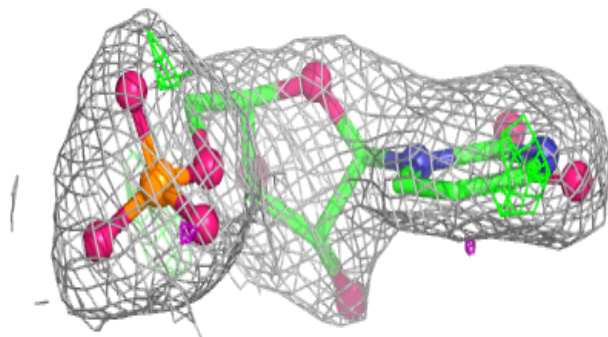
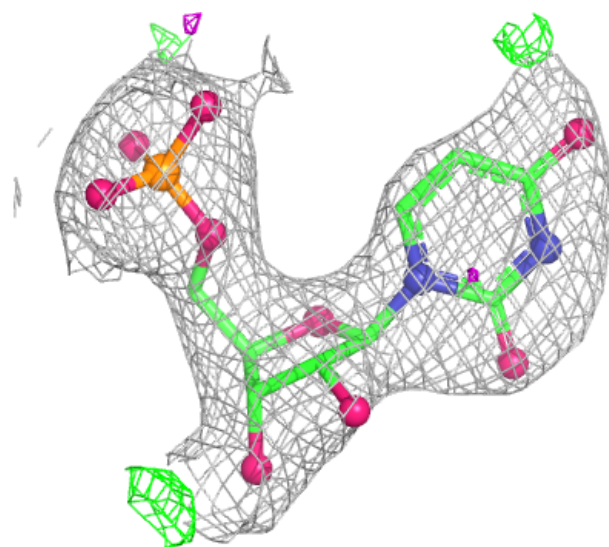
Electron density around U5P A 465:

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



Electron density around U5P B 456:

$2mF_o-DF_c$ (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.