



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:44 AM UTC

PDB ID : 6LD9 / pdb_00006ld9
Title : Crystal structure of cystathionine gamma synthase from *Xanthomonas oryzae* pv. *oryzae* in complex with cystathionine
Authors : Ngo, H.P.T.; Nguyen, T.D.Q.; Kang, L.W.
Deposited on : 2019-11-20
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

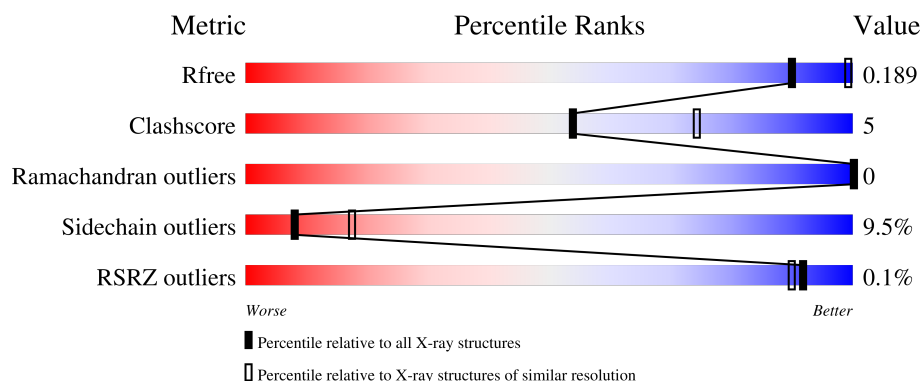
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	408	
1	B	408	
1	C	408	
1	D	408	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12052 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.

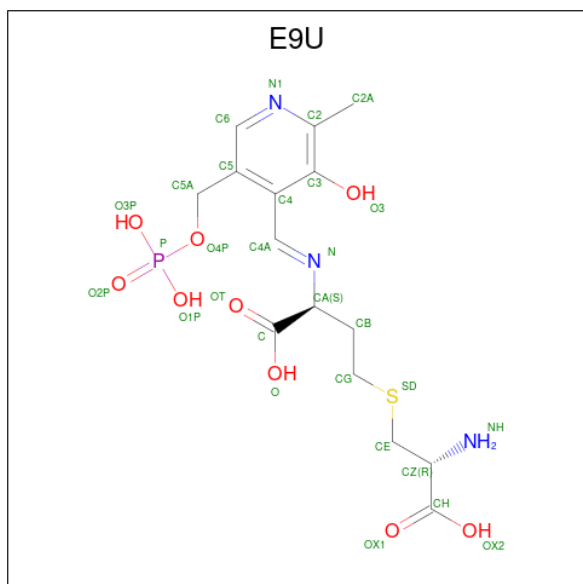
- Molecule 1 is a protein called Cystathionine gamma-synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	S	0	0	0
			2871	1813	506	545	7			
1	B	387	Total	C	N	O	S	0	0	0
			2871	1813	506	545	7			
1	C	388	Total	C	N	O	S	0	0	0
			2876	1816	507	546	7			
1	D	387	Total	C	N	O	S	0	0	0
			2871	1813	506	545	7			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q5H1U9
A	-1	SER	-	expression tag	UNP Q5H1U9
A	0	HIS	-	expression tag	UNP Q5H1U9
A	5	TYR	ASP	engineered mutation	UNP Q5H1U9
A	8	ASP	HIS	engineered mutation	UNP Q5H1U9
B	-2	GLY	-	expression tag	UNP Q5H1U9
B	-1	SER	-	expression tag	UNP Q5H1U9
B	0	HIS	-	expression tag	UNP Q5H1U9
B	5	TYR	ASP	engineered mutation	UNP Q5H1U9
B	8	ASP	HIS	engineered mutation	UNP Q5H1U9
C	-2	GLY	-	expression tag	UNP Q5H1U9
C	-1	SER	-	expression tag	UNP Q5H1U9
C	0	HIS	-	expression tag	UNP Q5H1U9
C	5	TYR	ASP	engineered mutation	UNP Q5H1U9
C	8	ASP	HIS	engineered mutation	UNP Q5H1U9
D	-2	GLY	-	expression tag	UNP Q5H1U9
D	-1	SER	-	expression tag	UNP Q5H1U9
D	0	HIS	-	expression tag	UNP Q5H1U9
D	5	TYR	ASP	engineered mutation	UNP Q5H1U9
D	8	ASP	HIS	engineered mutation	UNP Q5H1U9

- Molecule 2 is (2 {S})-4-[(2 {R})-2-azanyl-3-oxidanyl-3-oxidanylidene-propyl]sulfanyl-2-[({E})-2-methyl-3-oxidanyl-5-(phosphonooxymethyl)pyridin-4-yl]methylideneamino]butanoic acid (CCD ID: E9U) (formula: C₁₅H₂₂N₃O₉PS) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			29	15	3	9	1	1		
2	B	1	Total	C	N	O	P	S	0	0
			29	15	3	9	1	1		
2	C	1	Total	C	N	O	P	S	0	0
			29	15	3	9	1	1		
2	D	1	Total	C	N	O	P	S	0	0
			29	15	3	9	1	1		

- Molecule 3 is water.

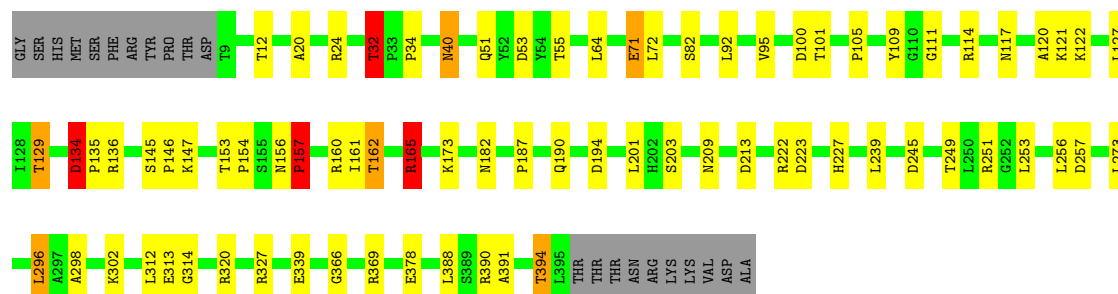
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	107	Total	O	0	0
			107	107		
3	B	104	Total	O	0	0
			104	104		
3	C	119	Total	O	0	0
			119	119		
3	D	117	Total	O	0	0
			117	117		

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.

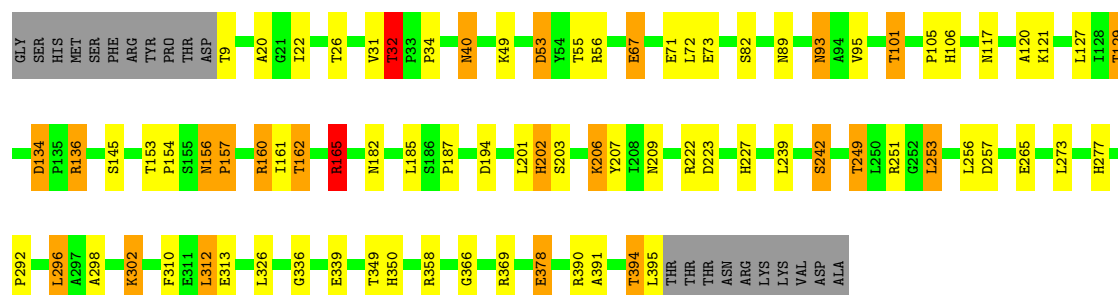
• Molecule 1: Cystathionine gamma-synthase

Chain A: 



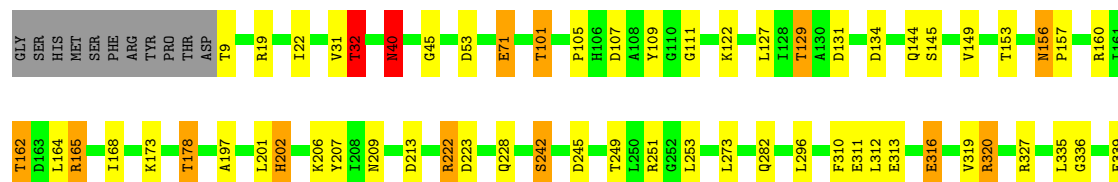
• Molecule 1: Cystathionine gamma-synthase

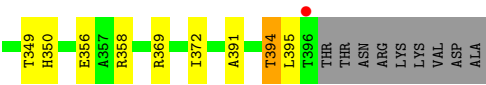
Chain B: 



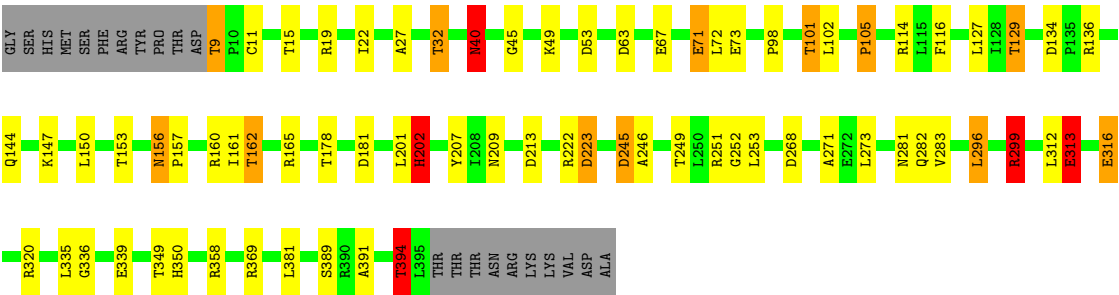
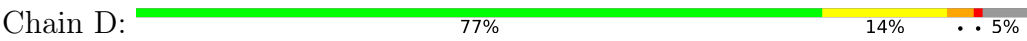
• Molecule 1: Cystathionine gamma-synthase

Chain C: 





● Molecule 1: Cystathionine gamma-synthase



4 Data and refinement statistics

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, α , β , γ	161.15Å 161.15Å 245.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.06 – 2.50 47.06 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.8 (47.06-2.50) 99.8 (47.06-2.50)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.50 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.153 , 0.186 0.157 , 0.189	Depositor DCC
R_{free} test set	5331 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	30.8	Xtriage
Anisotropy	0.114	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 9.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.478 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	12052	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.15% 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: E9U

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	1.41	15/2925 (0.5%)	1.63	23/3981 (0.6%)
1	B	1.38	21/2925 (0.7%)	1.61	24/3981 (0.6%)
1	C	1.38	13/2930 (0.4%)	1.57	21/3988 (0.5%)
1	D	1.39	15/2925 (0.5%)	1.57	19/3981 (0.5%)
All	All	1.39	64/11705 (0.5%)	1.60	87/15931 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	C	0	2
1	D	0	2
All	All	0	5

All (64) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	366	GLY	C-O	11.86	1.32	1.24
1	A	145	SER	CA-CB	-8.41	1.46	1.52
1	D	202	HIS	CE1-NE2	7.96	1.40	1.32
1	A	95	VAL	C-O	-7.76	1.16	1.24
1	A	105	PRO	C-O	-7.72	1.15	1.23
1	B	203	SER	CA-CB	-7.51	1.43	1.52
1	C	105	PRO	C-O	-7.41	1.15	1.23
1	C	202	HIS	CE1-NE2	7.33	1.39	1.32
1	B	93	ASN	C-O	-7.18	1.15	1.24
1	A	257	ASP	CG-OD2	7.13	1.39	1.25
1	A	51	GLN	C-O	-7.12	1.15	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	222	ARG	C-O	-7.12	1.15	1.24
1	B	257	ASP	CG-OD2	7.10	1.38	1.25
1	D	358	ARG	C-O	7.07	1.32	1.24
1	C	310	PHE	C-O	6.90	1.32	1.23
1	A	146	PRO	C-O	-6.85	1.16	1.23
1	A	20	ALA	C-O	6.80	1.31	1.24
1	B	358	ARG	C-O	6.71	1.31	1.24
1	D	71	GLU	C-O	-6.54	1.16	1.24
1	D	246	ALA	C-O	-6.52	1.16	1.24
1	B	145	SER	CA-CB	-6.51	1.48	1.52
1	B	34	PRO	C-O	-6.47	1.16	1.23
1	C	358	ARG	C-O	6.22	1.31	1.24
1	B	95	VAL	C-O	-6.22	1.18	1.24
1	B	277	HIS	CE1-NE2	6.17	1.38	1.32
1	C	71	GLU	C-O	-6.14	1.17	1.24
1	A	187	PRO	C-O	-6.12	1.16	1.24
1	D	381	LEU	C-O	-6.03	1.17	1.24
1	D	283	VAL	C-O	-6.02	1.17	1.24
1	D	40	ASN	C-O	6.01	1.31	1.23
1	C	31	VAL	C-O	-6.00	1.17	1.24
1	A	34	PRO	C-O	-5.96	1.17	1.23
1	C	149	VAL	C-O	5.96	1.30	1.24
1	B	89	ASN	C-O	-5.95	1.17	1.24
1	B	105	PRO	C-O	-5.94	1.17	1.23
1	D	45	GLY	C-O	5.93	1.32	1.23
1	B	366	GLY	C-O	5.91	1.30	1.23
1	A	314	GLY	C-O	5.90	1.32	1.24
1	A	92	LEU	C-O	-5.84	1.17	1.24
1	C	22	ILE	C-O	-5.83	1.17	1.24
1	B	22	ILE	C-O	-5.80	1.17	1.24
1	B	292	PRO	C-O	-5.80	1.16	1.24
1	A	203	SER	CA-CB	-5.59	1.46	1.53
1	C	40	ASN	C-O	5.53	1.30	1.23
1	B	187	PRO	C-O	-5.52	1.17	1.24
1	B	253	LEU	C-O	-5.44	1.17	1.24
1	C	245	ASP	CG-OD2	5.43	1.35	1.25
1	B	20	ALA	C-O	5.42	1.30	1.24
1	D	268	ASP	C-O	-5.38	1.17	1.24
1	A	298	ALA	C-O	-5.37	1.17	1.24
1	D	22	ILE	C-O	-5.37	1.17	1.24
1	B	121	LYS	C-O	-5.36	1.17	1.24
1	C	222	ARG	C-O	-5.30	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	129	THR	C-O	-5.28	1.17	1.24
1	B	249	THR	C-O	-5.27	1.18	1.24
1	B	26	THR	C-O	-5.25	1.17	1.24
1	D	245	ASP	CG-OD2	5.21	1.35	1.25
1	D	105	PRO	C-O	-5.17	1.18	1.23
1	B	298	ALA	C-O	-5.12	1.17	1.24
1	C	45	GLY	C-O	5.09	1.30	1.23
1	D	102	LEU	C-O	5.09	1.30	1.24
1	C	131	ASP	C-O	-5.07	1.17	1.23
1	B	106	HIS	CE1-NE2	5.04	1.37	1.32
1	D	27	ALA	C-O	-5.02	1.18	1.24

All (87) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	129	THR	CB-CA-C	10.60	128.07	110.79
1	A	134	ASP	CB-CA-C	10.01	121.28	110.17
1	B	32	THR	CA-CB-OG1	-10.00	94.60	109.60
1	A	32	THR	CA-CB-OG1	-9.97	94.64	109.60
1	A	114	ARG	CG-CD-NE	9.67	133.28	112.00
1	D	134	ASP	CA-CB-CG	-9.43	103.17	112.60
1	B	129	THR	CB-CA-C	8.90	125.30	110.79
1	D	32	THR	CA-CB-OG1	-8.83	96.35	109.60
1	B	134	ASP	CB-CA-C	8.35	119.44	110.17
1	A	53	ASP	CA-CB-CG	8.19	120.78	112.60
1	C	32	THR	CA-CB-OG1	-7.61	98.19	109.60
1	A	134	ASP	CA-CB-CG	-7.51	105.09	112.60
1	A	339	GLU	CB-CG-CD	7.43	125.24	112.60
1	C	134	ASP	N-CA-CB	-7.11	99.30	111.39
1	C	356	GLU	CB-CG-CD	7.03	124.54	112.60
1	C	165	ARG	CG-CD-NE	-6.95	96.72	112.00
1	C	134	ASP	CA-CB-CG	-6.91	105.69	112.60
1	C	165	ARG	CB-CG-CD	6.88	127.12	111.30
1	A	388	LEU	CA-C-O	-6.80	113.34	120.55
1	D	129	THR	CB-CA-C	6.75	121.49	110.22
1	B	227	HIS	CA-C-O	-6.74	113.74	120.82
1	D	223	ASP	CB-CA-C	6.69	120.72	109.48
1	C	165	ARG	CB-CA-C	-6.64	100.45	110.88
1	C	164	LEU	CA-C-O	-6.59	113.90	120.82
1	C	129	THR	CB-CA-C	6.54	121.25	110.78
1	D	160	ARG	CG-CD-NE	-6.53	97.63	112.00
1	D	67	GLU	N-CA-C	-6.53	104.24	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	117	ASN	CB-CA-C	6.48	121.87	110.85
1	B	117	ASN	CB-CA-C	6.45	123.54	110.38
1	D	101	THR	CB-CA-C	6.45	120.98	109.38
1	C	134	ASP	CB-CA-C	6.30	119.70	111.01
1	B	194	ASP	CA-CB-CG	6.29	118.89	112.60
1	B	120	ALA	CA-C-O	-6.23	113.81	120.42
1	B	129	THR	CA-CB-OG1	-6.17	100.34	109.60
1	D	394	THR	CA-CB-OG1	-6.05	100.52	109.60
1	B	134	ASP	CA-CB-CG	-6.04	106.56	112.60
1	A	227	HIS	CA-C-O	-6.00	114.52	120.82
1	D	313	GLU	CB-CG-CD	6.00	122.80	112.60
1	C	101	THR	CB-CA-C	5.94	120.08	109.38
1	C	206	LYS	CB-CA-C	5.90	122.16	110.42
1	C	160	ARG	CG-CD-NE	-5.87	99.08	112.00
1	D	134	ASP	N-CA-CB	-5.85	101.78	111.78
1	B	67	GLU	CB-CA-C	5.82	120.75	110.85
1	B	53	ASP	CA-CB-CG	5.80	118.40	112.60
1	A	100	ASP	CA-CB-CG	5.78	118.38	112.60
1	A	12	THR	CA-C-N	5.66	128.18	120.54
1	A	12	THR	C-N-CA	5.66	128.18	120.54
1	B	265	GLU	CB-CA-C	5.65	120.17	110.79
1	A	71	GLU	CB-CG-CD	5.61	122.14	112.60
1	B	339	GLU	CB-CG-CD	5.60	122.11	112.60
1	D	339	GLU	CB-CG-CD	5.59	122.11	112.60
1	D	116	PHE	CA-C-O	-5.56	114.65	120.55
1	A	194	ASP	CA-CB-CG	5.53	118.13	112.60
1	D	271	ALA	CA-C-O	-5.53	115.01	120.82
1	C	395	LEU	CA-C-N	5.49	131.59	121.70
1	C	395	LEU	C-N-CA	5.49	131.59	121.70
1	D	53	ASP	CA-CB-CG	5.47	118.07	112.60
1	B	160	ARG	CB-CA-C	5.45	119.57	109.54
1	D	389	SER	CB-CA-C	-5.43	102.35	110.88
1	C	339	GLU	CB-CG-CD	5.43	121.83	112.60
1	A	165	ARG	NE-CZ-NH2	5.42	124.08	119.20
1	D	299	ARG	CB-CG-CD	5.41	123.74	111.30
1	B	206	LYS	CB-CA-C	5.37	121.11	110.42
1	B	378	GLU	CB-CG-CD	5.36	121.71	112.60
1	C	53	ASP	CA-CB-CG	5.33	117.93	112.60
1	B	390	ARG	N-CA-C	-5.33	105.55	111.36
1	A	120	ALA	CA-C-O	-5.29	114.81	120.42
1	B	302	LYS	N-CA-C	-5.29	106.33	112.89
1	C	107	ASP	CA-CB-CG	5.28	117.88	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	390	ARG	N-CA-C	-5.28	105.52	111.28
1	D	63	ASP	CA-CB-CG	5.28	117.88	112.60
1	B	136	ARG	CA-C-N	5.26	127.64	120.54
1	B	136	ARG	C-N-CA	5.26	127.64	120.54
1	A	114	ARG	NE-CZ-NH2	-5.25	114.47	119.20
1	A	256	LEU	CA-C-O	-5.24	115.31	120.82
1	B	165	ARG	CG-CD-NE	-5.24	100.48	112.00
1	A	157	PRO	N-CA-CB	-5.21	96.87	102.60
1	B	67	GLU	N-CA-C	-5.19	105.71	111.36
1	C	206	LYS	CG-CD-CE	5.19	123.23	111.30
1	B	256	LEU	CA-C-O	-5.17	115.39	120.82
1	C	319	VAL	CA-C-O	-5.15	115.91	121.27
1	D	389	SER	N-CA-CB	5.14	117.47	110.01
1	A	64	LEU	N-CA-C	-5.14	105.76	111.36
1	D	252	GLY	O-C-N	5.10	127.08	122.18
1	C	372	ILE	CB-CA-C	5.10	117.13	110.96
1	A	51	GLN	CA-C-O	-5.02	115.53	120.90
1	B	326	LEU	CA-C-O	-5.00	115.22	120.92

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	202	HIS	Peptide
1	C	202	HIS	Peptide
1	C	9	THR	Peptide
1	D	202	HIS	Peptide
1	D	9	THR	Peptide

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	2871	0	2868	31	0
1	B	2871	0	2868	41	0
1	C	2876	0	2870	28	0
1	D	2871	0	2868	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	29	0	0	0	0
2	B	29	0	0	2	0
2	C	29	0	0	1	0
2	D	29	0	0	3	0
3	A	107	0	0	1	0
3	B	104	0	0	7	0
3	C	119	0	0	5	0
3	D	117	0	0	3	0
All	All	12052	0	11474	113	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 5.

All (113) 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:A:251:ARG:HH22	1:C:209:ASN:HD21	1.32	0.77
1:A:165:ARG:HG2	1:A:165:ARG:HH21	1.52	0.72
1:B:40:ASN:H	1:B:40:ASN:HD22	1.39	0.71
1:A:201:LEU:HD12	1:A:201:LEU:C	2.17	0.69
1:D:316:GLU:OE2	1:D:320:ARG:NH2	2.25	0.69
1:B:209:ASN:HD21	1:D:251:ARG:HH22	1.41	0.69
1:B:391:ALA:O	1:B:394:THR:HB	1.92	0.69
1:A:209:ASN:HD21	1:C:251:ARG:HH22	1.41	0.69
1:A:134:ASP:OD1	1:A:134:ASP:C	2.34	0.68
1:C:156:ASN:HD21	1:C:369:ARG:HH11	1.40	0.68
1:A:40:ASN:HD22	1:A:40:ASN:H	1.41	0.68
1:B:162:THR:HG22	3:B:948:HOH:O	1.94	0.67
1:B:134:ASP:OD1	1:B:134:ASP:C	2.37	0.67
1:B:251:ARG:HH22	1:D:209:ASN:HD21	1.43	0.67
1:B:209:ASN:HD21	1:D:251:ARG:NH2	1.94	0.66
1:A:209:ASN:HD22	1:A:249:THR:HA	1.62	0.65
1:B:201:LEU:C	1:B:201:LEU:HD12	2.22	0.64
1:C:153:THR:HB	1:C:162:THR:HG23	1.81	0.63
1:A:251:ARG:NH2	1:C:209:ASN:HD21	1.96	0.63
1:C:209:ASN:HD22	1:C:249:THR:HA	1.64	0.62
1:B:101:THR:HG23	3:B:902:HOH:O	1.99	0.62
1:B:40:ASN:HD22	1:B:40:ASN:N	1.96	0.61
1:D:349:THR:OG1	1:D:350:HIS:HD2	1.84	0.61
1:A:153:THR:HB	1:A:162:THR:HG23	1.81	0.61
1:D:153:THR:HB	1:D:162:THR:HG23	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:156:ASN:HD21	1:D:369:ARG:HH11	1.50	0.60
1:A:209:ASN:HD21	1:C:251:ARG:NH2	1.99	0.59
1:B:209:ASN:HD22	1:B:249:THR:HA	1.67	0.59
1:D:391:ALA:O	1:D:394:THR:HB	2.03	0.59
1:A:40:ASN:HD22	1:A:40:ASN:N	1.98	0.58
1:B:156:ASN:HD21	1:B:369:ARG:HH11	1.51	0.58
1:B:251:ARG:NH2	1:D:209:ASN:HD21	2.02	0.57
1:D:201:LEU:HD12	1:D:201:LEU:C	2.30	0.57
1:D:162:THR:HG22	3:D:943:HOH:O	2.05	0.57
1:A:156:ASN:HD21	1:A:369:ARG:HH11	1.52	0.56
1:D:162:THR:CG2	3:D:943:HOH:O	2.51	0.56
1:B:162:THR:CG2	3:B:948:HOH:O	2.53	0.56
1:C:197:ALA:O	1:C:222:ARG:NH2	2.39	0.56
1:C:349:THR:OG1	1:C:350:HIS:HD2	1.88	0.56
1:A:391:ALA:O	1:A:394:THR:HB	2.05	0.55
1:A:32:THR:HG21	1:C:213:ASP:HB3	1.89	0.55
1:B:223:ASP:HB3	3:B:975:HOH:O	2.06	0.54
1:D:40:ASN:HD22	1:D:40:ASN:H	1.54	0.54
1:A:162:THR:HG21	3:A:966:HOH:O	2.08	0.53
1:C:109:TYR:CE2	1:C:111:GLY:HA3	2.44	0.53
1:D:209:ASN:HD22	1:D:249:THR:HA	1.73	0.53
1:C:162:THR:HG22	3:C:948:HOH:O	2.08	0.53
1:C:391:ALA:O	1:C:394:THR:HB	2.07	0.53
1:B:153:THR:HB	1:B:162:THR:HG23	1.91	0.52
1:A:122:LYS:NZ	1:B:93:ASN:O	2.42	0.52
1:C:19:ARG:HD3	3:C:943:HOH:O	2.09	0.52
1:B:206:LYS:HZ1	2:B:801:E9U:C4A	2.23	0.51
1:A:109:TYR:CE2	1:A:111:GLY:HA3	2.45	0.51
1:B:32:THR:HG21	1:D:213:ASP:HB3	1.92	0.51
1:C:282:GLN:NE2	3:C:904:HOH:O	2.44	0.51
1:B:161:ILE:HG22	1:B:296:LEU:HD13	1.93	0.50
1:C:162:THR:HG21	3:C:1006:HOH:O	2.10	0.50
1:B:310:PHE:HE2	1:B:312:LEU:HD13	1.75	0.50
1:C:40:ASN:HD22	1:C:40:ASN:H	1.58	0.50
1:D:73:GLU:OE2	1:D:202:HIS:NE2	2.37	0.49
1:B:310:PHE:CE2	1:B:312:LEU:HD13	2.46	0.49
1:B:162:THR:HG21	3:B:967:HOH:O	2.13	0.47
1:A:156:ASN:HD22	1:A:157:PRO:CA	2.27	0.47
1:C:168:ILE:HG12	1:C:178:THR:HG21	1.97	0.47
1:B:165:ARG:HD2	3:B:977:HOH:O	2.13	0.47
1:B:49:LYS:HD2	1:B:53:ASP:OD2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:162:THR:CG2	3:C:948:HOH:O	2.62	0.46
1:B:156:ASN:HA	1:B:157:PRO:HA	1.75	0.46
1:B:349:THR:OG1	1:B:350:HIS:HD2	1.98	0.46
1:C:122:LYS:O	1:D:98:PRO:HD2	2.14	0.46
1:B:207:TYR:CE1	1:B:336:GLY:HA2	2.51	0.46
1:D:11:CYS:HB2	1:D:15:THR:HB	1.98	0.46
1:D:161:ILE:HG22	1:D:296:LEU:HD13	1.97	0.46
1:B:40:ASN:N	1:B:40:ASN:ND2	2.64	0.45
1:D:114:ARG:HH22	2:D:801:E9U:CH	2.29	0.45
1:D:40:ASN:HD22	1:D:40:ASN:N	2.13	0.44
2:B:801:E9U:O3	2:B:801:E9U:N	2.50	0.44
1:B:101:THR:CG2	3:B:902:HOH:O	2.61	0.44
1:C:201:LEU:C	1:C:201:LEU:HD12	2.43	0.44
1:A:251:ARG:HH22	1:C:209:ASN:ND2	2.08	0.44
1:A:201:LEU:C	1:A:201:LEU:CD1	2.90	0.43
1:C:316:GLU:OE2	1:C:320:ARG:NH2	2.51	0.43
1:A:182:ASN:HD21	1:A:190:GLN:HB3	1.83	0.43
1:C:207:TYR:CE1	1:C:336:GLY:HA2	2.53	0.43
1:D:19:ARG:HD3	3:D:981:HOH:O	2.17	0.43
1:A:153:THR:HA	1:A:154:PRO:C	2.40	0.43
1:D:281:ASN:HB2	1:D:313:GLU:OE1	2.18	0.43
1:B:154:PRO:HD3	1:B:185:LEU:CD1	2.48	0.43
1:D:181:ASP:OD2	2:D:801:E9U:N1	2.51	0.43
1:B:156:ASN:HD22	1:B:157:PRO:CA	2.32	0.43
2:C:801:E9U:N	2:C:801:E9U:O3	2.52	0.43
1:B:209:ASN:ND2	1:D:251:ARG:HH22	2.11	0.42
1:C:242:SER:HB2	1:D:245:ASP:OD2	2.18	0.42
1:A:82:SER:HB3	1:B:82:SER:HB3	2.00	0.42
1:C:40:ASN:HD22	1:C:40:ASN:N	2.15	0.42
1:D:207:TYR:CE1	1:D:336:GLY:HA2	2.55	0.42
1:A:245:ASP:OD2	1:B:242:SER:HB2	2.19	0.42
1:B:31:VAL:HB	1:C:40:ASN:ND2	2.35	0.42
2:D:801:E9U:O3	2:D:801:E9U:N	2.53	0.41
1:A:40:ASN:N	1:A:40:ASN:ND2	2.65	0.41
1:B:73:GLU:OE2	1:B:202:HIS:NE2	2.48	0.41
1:A:134:ASP:HA	1:A:135:PRO:HD3	1.99	0.41
1:A:121:LYS:HE2	1:A:121:LYS:HB3	1.85	0.41
1:D:165:ARG:HG2	1:D:165:ARG:NH1	2.36	0.41
1:D:299:ARG:HB2	1:D:299:ARG:NH1	2.36	0.41
1:A:213:ASP:HB3	1:C:32:THR:HG21	2.03	0.41
1:D:105:PRO:HD3	1:D:150:LEU:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:56:ARG:HH11	1:B:56:ARG:HG2	1.85	0.40
1:B:153:THR:HA	1:B:154:PRO:C	2.46	0.40
1:A:182:ASN:HD22	1:A:182:ASN:HA	1.74	0.40
1:A:161:ILE:HG22	1:A:296:LEU:HD13	2.02	0.40
1:A:239:LEU:HD23	1:B:239:LEU:HD23	2.02	0.40
1:B:182:ASN:HD22	1:B:182:ASN:HA	1.72	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	385/408 (94%)	372 (97%)	13 (3%)	0	100	100
1	B	385/408 (94%)	371 (96%)	14 (4%)	0	100	100
1	C	386/408 (95%)	376 (97%)	10 (3%)	0	100	100
1	D	385/408 (94%)	375 (97%)	10 (3%)	0	100	100
All	All	1541/1632 (94%)	1494 (97%)	47 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	291/310 (94%)	262 (90%)	29 (10%)	7	15
1	B	291/310 (94%)	264 (91%)	27 (9%)	8	18
1	C	291/310 (94%)	263 (90%)	28 (10%)	8	17
1	D	291/310 (94%)	264 (91%)	27 (9%)	8	18
All	All	1164/1240 (94%)	1053 (90%)	111 (10%)	8	17

All (111) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	ARG
1	A	32	THR
1	A	40	ASN
1	A	55	THR
1	A	71	GLU
1	A	72	LEU
1	A	101	THR
1	A	127	LEU
1	A	129	THR
1	A	134	ASP
1	A	136	ARG
1	A	147	LYS
1	A	157	PRO
1	A	160	ARG
1	A	162	THR
1	A	165	ARG
1	A	173	LYS
1	A	222	ARG
1	A	223	ASP
1	A	253	LEU
1	A	273	LEU
1	A	296	LEU
1	A	302	LYS
1	A	312	LEU
1	A	313	GLU
1	A	320	ARG
1	A	327	ARG
1	A	378	GLU
1	A	394	THR
1	B	9	THR
1	B	32	THR
1	B	40	ASN
1	B	55	THR

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Mol	Chain	Res	Type
1	B	67	GLU
1	B	71	GLU
1	B	72	LEU
1	B	101	THR
1	B	127	LEU
1	B	129	THR
1	B	136	ARG
1	B	156	ASN
1	B	157	PRO
1	B	160	ARG
1	B	162	THR
1	B	165	ARG
1	B	222	ARG
1	B	242	SER
1	B	253	LEU
1	B	273	LEU
1	B	296	LEU
1	B	302	LYS
1	B	312	LEU
1	B	313	GLU
1	B	378	GLU
1	B	394	THR
1	B	395	LEU
1	C	32	THR
1	C	40	ASN
1	C	71	GLU
1	C	101	THR
1	C	127	LEU
1	C	129	THR
1	C	144	GLN
1	C	145	SER
1	C	156	ASN
1	C	157	PRO
1	C	162	THR
1	C	165	ARG
1	C	173	LYS
1	C	178	THR
1	C	223	ASP
1	C	228	GLN
1	C	242	SER
1	C	253	LEU
1	C	273	LEU

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Mol	Chain	Res	Type
1	C	296	LEU
1	C	311	GLU
1	C	312	LEU
1	C	313	GLU
1	C	316	GLU
1	C	320	ARG
1	C	327	ARG
1	C	335	LEU
1	C	394	THR
1	D	9	THR
1	D	32	THR
1	D	40	ASN
1	D	49	LYS
1	D	71	GLU
1	D	72	LEU
1	D	101	THR
1	D	127	LEU
1	D	129	THR
1	D	136	ARG
1	D	144	GLN
1	D	147	LYS
1	D	156	ASN
1	D	157	PRO
1	D	162	THR
1	D	178	THR
1	D	223	ASP
1	D	253	LEU
1	D	273	LEU
1	D	282	GLN
1	D	296	LEU
1	D	299	ARG
1	D	312	LEU
1	D	313	GLU
1	D	316	GLU
1	D	335	LEU
1	D	394	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	48	ASN

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Mol	Chain	Res	Type
1	A	117	ASN
1	A	156	ASN
1	A	182	ASN
1	A	209	ASN
1	A	281	ASN
1	A	282	GLN
1	A	301	GLN
1	A	350	HIS
1	B	40	ASN
1	B	48	ASN
1	B	97	GLN
1	B	117	ASN
1	B	156	ASN
1	B	182	ASN
1	B	209	ASN
1	B	301	GLN
1	B	350	HIS
1	C	40	ASN
1	C	97	GLN
1	C	117	ASN
1	C	156	ASN
1	C	182	ASN
1	C	209	ASN
1	C	301	GLN
1	C	350	HIS
1	D	40	ASN
1	D	97	GLN
1	D	117	ASN
1	D	156	ASN
1	D	182	ASN
1	D	209	ASN
1	D	227	HIS
1	D	301	GLN
1	D	350	HIS

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	E9U	A	801	-	28,29,29	1.72	7 (25%)	33,40,40	2.83	10 (30%)
2	E9U	B	801	-	28,29,29	1.95	8 (28%)	33,40,40	2.67	9 (27%)
2	E9U	D	801	-	28,29,29	1.97	6 (21%)	33,40,40	2.29	12 (36%)
2	E9U	C	801	-	28,29,29	1.92	7 (25%)	33,40,40	2.12	11 (33%)

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	E9U	A	801	-	-	6/26/26/26	0/1/1/1
2	E9U	B	801	-	-	6/26/26/26	0/1/1/1
2	E9U	D	801	-	-	10/26/26/26	0/1/1/1
2	E9U	C	801	-	-	13/26/26/26	0/1/1/1

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	801	E9U	C3-C2	6.83	1.48	1.41
2	C	801	E9U	C3-C2	5.44	1.46	1.41
2	B	801	E9U	C3-C2	5.29	1.46	1.41
2	A	801	E9U	C3-C2	4.78	1.45	1.41
2	B	801	E9U	CB-CA	4.52	1.58	1.53
2	C	801	E9U	C4-C5	4.44	1.48	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	E9U	OT-C	3.82	1.33	1.22
2	B	801	E9U	CA-N	3.63	1.50	1.46
2	D	801	E9U	C4-C5	3.59	1.47	1.42
2	D	801	E9U	OX2-CH	-3.36	1.19	1.30
2	B	801	E9U	C4-C3	3.33	1.46	1.41
2	D	801	E9U	C4-C3	3.21	1.46	1.41
2	A	801	E9U	C4-C3	2.92	1.45	1.41
2	A	801	E9U	OX2-CH	-2.88	1.21	1.30
2	B	801	E9U	C4-C5	2.86	1.45	1.42
2	A	801	E9U	CB-CA	2.69	1.56	1.53
2	C	801	E9U	C2-N1	2.66	1.38	1.33
2	A	801	E9U	C4-C5	2.60	1.45	1.42
2	C	801	E9U	OX2-CH	-2.52	1.22	1.30
2	C	801	E9U	C4-C3	2.45	1.45	1.41
2	B	801	E9U	OT-C	2.41	1.29	1.22
2	D	801	E9U	OT-C	2.40	1.29	1.22
2	A	801	E9U	CA-N	2.37	1.49	1.46
2	B	801	E9U	C6-C5	2.25	1.42	1.37
2	B	801	E9U	OX2-CH	-2.21	1.23	1.30
2	C	801	E9U	P-O4P	2.03	1.66	1.60
2	C	801	E9U	C6-C5	2.02	1.41	1.37
2	D	801	E9U	C6-C5	2.00	1.41	1.37

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	E9U	CA-N-C4A	6.99	127.15	116.77
2	B	801	E9U	C2A-C2-C3	-6.65	113.02	120.80
2	A	801	E9U	CE-SD-CG	6.64	121.97	102.26
2	A	801	E9U	C2A-C2-C3	-6.56	113.12	120.80
2	B	801	E9U	CE-SD-CG	6.27	120.87	102.26
2	A	801	E9U	C3-C4-C5	-5.90	113.54	118.28
2	B	801	E9U	CA-N-C4A	5.42	124.82	116.77
2	C	801	E9U	CZ-CE-SD	5.24	132.84	113.51
2	D	801	E9U	CZ-CE-SD	5.07	132.22	113.51
2	B	801	E9U	C3-C4-C5	-4.97	114.29	118.28
2	D	801	E9U	C4-C3-C2	-4.84	117.42	120.14
2	A	801	E9U	OX2-CH-OX1	-4.48	113.91	124.08
2	D	801	E9U	C3-C4-C5	-4.44	114.71	118.28
2	B	801	E9U	OX2-CH-OX1	-4.40	114.11	124.08
2	B	801	E9U	C5-C4-C4A	4.35	128.18	121.47
2	C	801	E9U	C4-C3-C2	-4.31	117.72	120.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	801	E9U	CA-N-C4A	4.25	123.09	116.77
2	C	801	E9U	CA-N-C4A	4.09	122.84	116.77
2	A	801	E9U	CB-CG-SD	-3.99	104.55	113.45
2	B	801	E9U	CB-CG-SD	-3.79	105.00	113.45
2	D	801	E9U	CE-SD-CG	3.51	112.67	102.26
2	A	801	E9U	C5-C4-C4A	3.50	126.86	121.47
2	B	801	E9U	C2A-C2-N1	3.49	124.22	117.64
2	C	801	E9U	CE-SD-CG	3.33	112.16	102.26
2	A	801	E9U	C2A-C2-N1	3.20	123.66	117.64
2	D	801	E9U	CB-CA-C	3.20	114.79	109.48
2	C	801	E9U	CB-CA-C	3.18	114.76	109.48
2	C	801	E9U	C3-C4-C5	-3.15	115.75	118.28
2	D	801	E9U	CE-CZ-CH	-2.99	102.12	110.70
2	D	801	E9U	C5-C4-C4A	2.84	125.86	121.47
2	C	801	E9U	OX2-CH-OX1	-2.79	117.75	124.08
2	D	801	E9U	O3-C3-C2	2.78	123.34	117.58
2	C	801	E9U	O4P-P-O2P	-2.70	99.13	106.44
2	B	801	E9U	O3P-P-O4P	2.62	113.51	106.67
2	C	801	E9U	C5-C4-C4A	2.60	125.48	121.47
2	D	801	E9U	OX2-CH-OX1	-2.60	118.18	124.08
2	C	801	E9U	C2A-C2-C3	-2.49	117.89	120.80
2	A	801	E9U	O1P-P-O4P	2.38	112.87	106.67
2	A	801	E9U	O-C-OT	2.33	129.36	124.08
2	C	801	E9U	O3-C3-C2	2.17	122.07	117.58
2	D	801	E9U	C3-C2-N1	2.07	123.56	120.96
2	D	801	E9U	C6-C5-C4	2.01	121.66	118.21

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	E9U	C-CA-N-C4A
2	A	801	E9U	CB-CA-N-C4A
2	A	801	E9U	N-CA-CB-CG
2	A	801	E9U	C-CA-CB-CG
2	A	801	E9U	SD-CE-CZ-CH
2	A	801	E9U	SD-CE-CZ-NH
2	B	801	E9U	CB-CA-N-C4A
2	B	801	E9U	N-CA-CB-CG
2	B	801	E9U	C-CA-CB-CG
2	B	801	E9U	SD-CE-CZ-CH
2	B	801	E9U	SD-CE-CZ-NH

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Mol	Chain	Res	Type	Atoms
2	C	801	E9U	C-CA-N-C4A
2	C	801	E9U	CB-CA-N-C4A
2	C	801	E9U	OT-C-CA-CB
2	C	801	E9U	N-CA-CB-CG
2	C	801	E9U	C-CA-CB-CG
2	C	801	E9U	SD-CE-CZ-CH
2	C	801	E9U	SD-CE-CZ-NH
2	D	801	E9U	C-CA-N-C4A
2	D	801	E9U	CB-CA-N-C4A
2	D	801	E9U	SD-CE-CZ-CH
2	D	801	E9U	SD-CE-CZ-NH
2	D	801	E9U	CA-CB-CG-SD
2	D	801	E9U	OX1-CH-CZ-CE
2	D	801	E9U	OX2-CH-CZ-CE
2	C	801	E9U	CZ-CE-SD-CG
2	C	801	E9U	CA-CB-CG-SD
2	C	801	E9U	O-C-CA-CB
2	D	801	E9U	OT-C-CA-CB
2	C	801	E9U	OT-C-CA-N
2	D	801	E9U	CZ-CE-SD-CG
2	D	801	E9U	O-C-CA-CB
2	C	801	E9U	OX1-CH-CZ-CE
2	C	801	E9U	OX2-CH-CZ-CE
2	B	801	E9U	C-CA-N-C4A

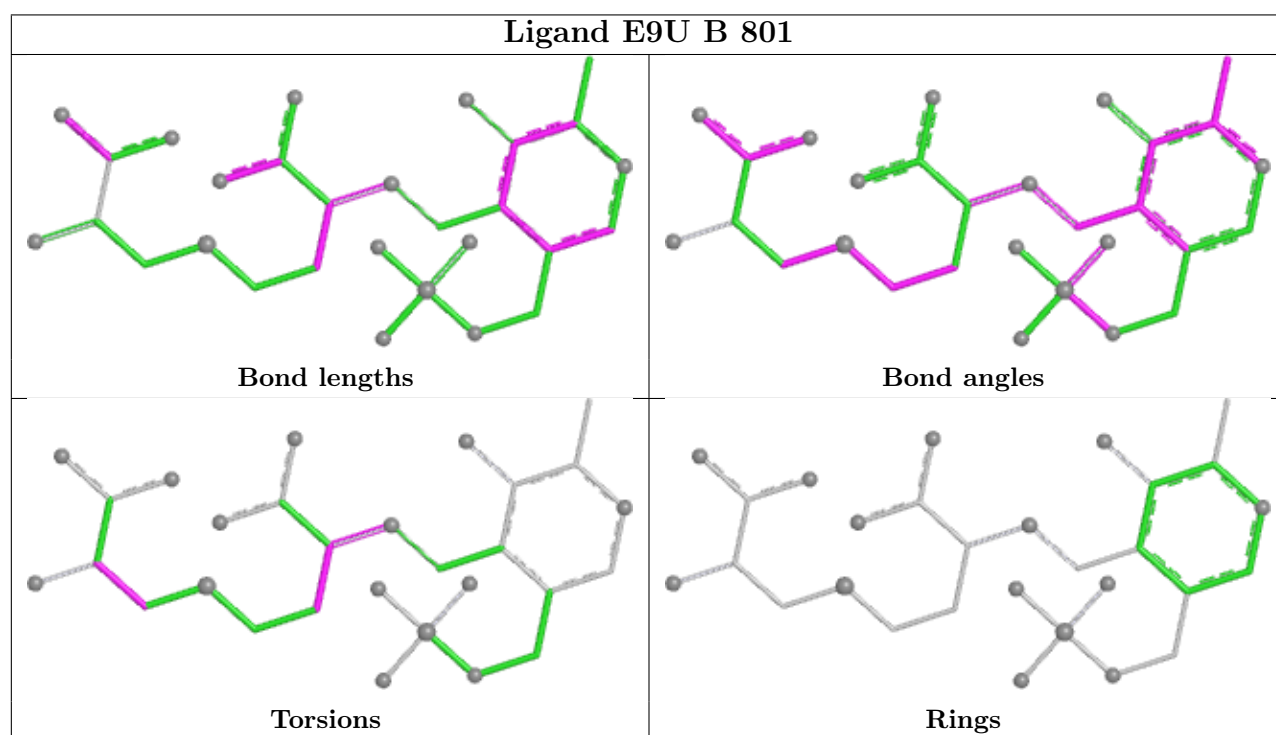
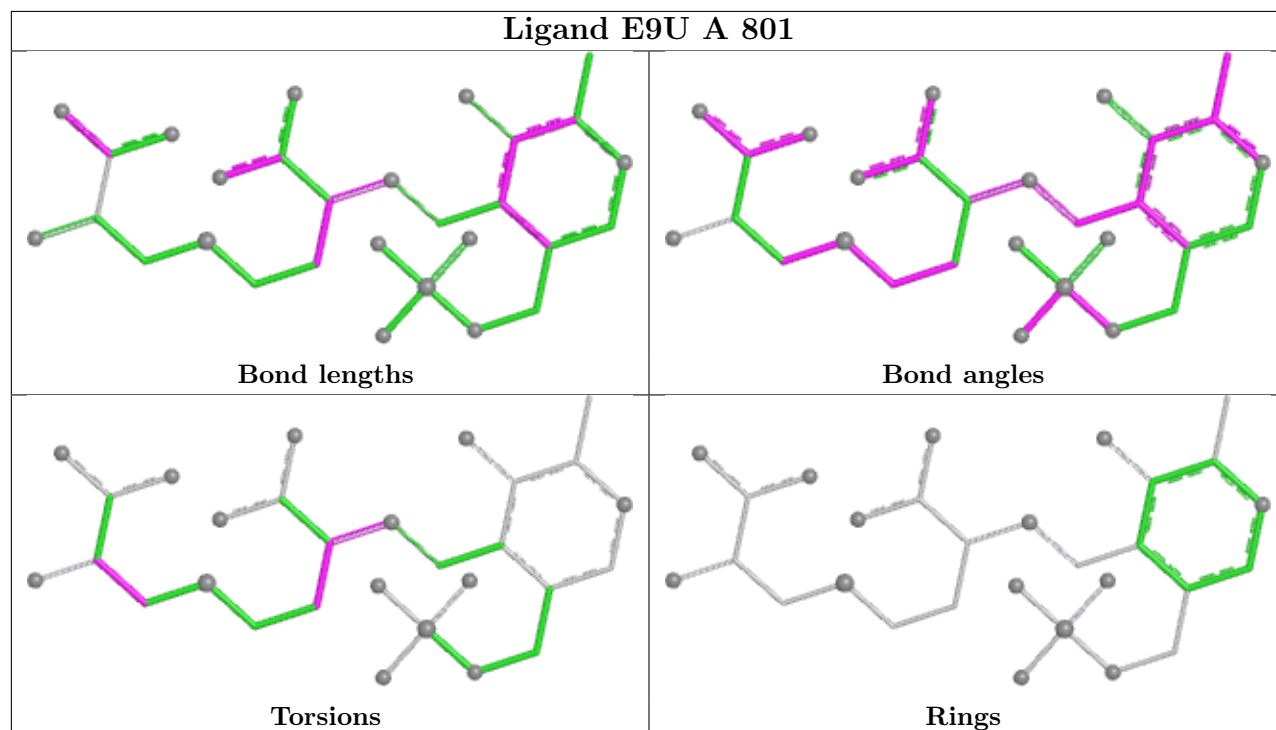
There are no ring outliers.

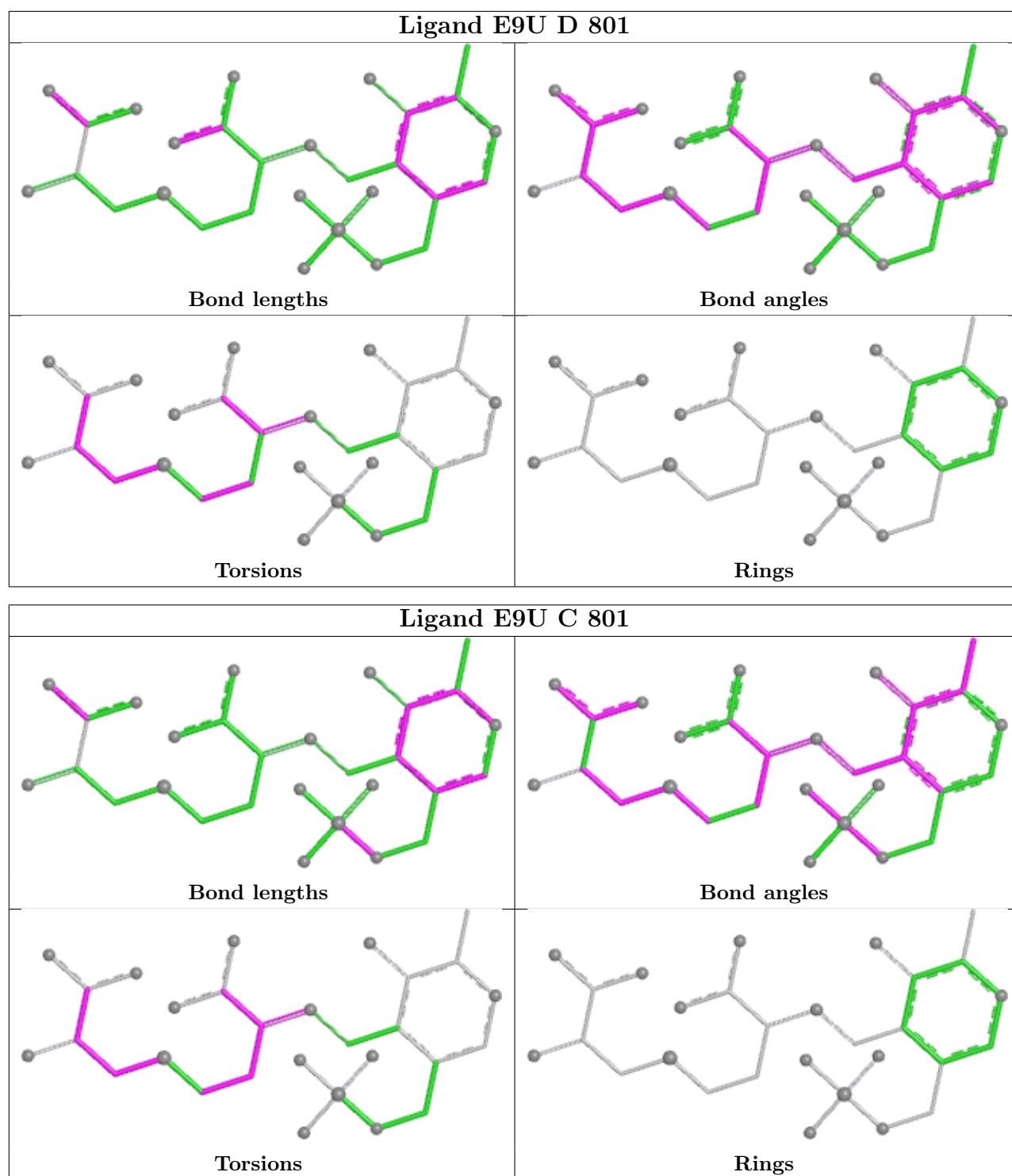
3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	801	E9U	2	0
2	D	801	E9U	3	0
2	C	801	E9U	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	A	387/408 (94%)	-1.72	0	100	100	18, 30, 52, 80	0
1	B	387/408 (94%)	-1.71	0	100	100	19, 30, 51, 85	0
1	C	388/408 (95%)	-1.68	1 (0%)	90	87	20, 31, 53, 102	0
1	D	387/408 (94%)	-1.70	0	100	100	19, 31, 54, 94	0
All	All	1549/1632 (94%)	-1.70	1 (0%)	92	90	18, 31, 54, 102	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	396	THR	2.5

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
2	E9U	A	801	29/29	1.00	0.02	22,29,43,56	0

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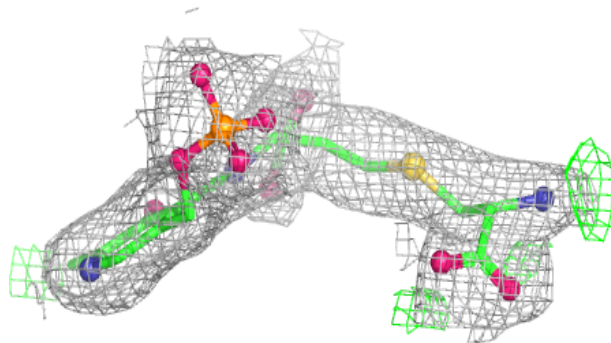
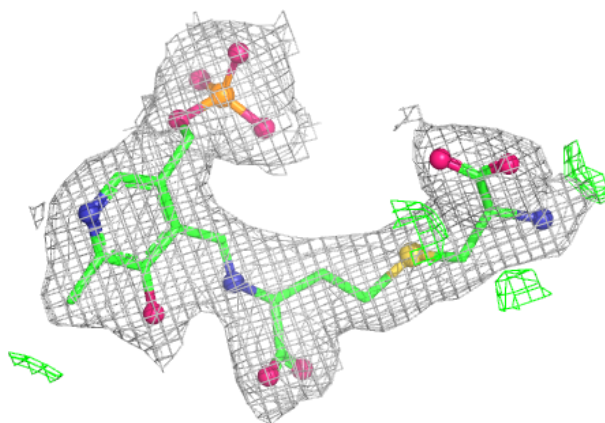
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	E9U	B	801	29/29	1.00	0.02	23,28,43,57	0
2	E9U	C	801	29/29	1.00	0.02	21,27,44,63	0
2	E9U	D	801	29/29	1.00	0.02	22,28,42,64	0

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.

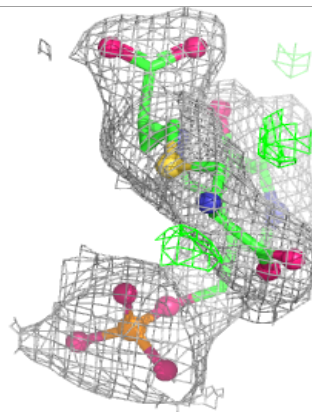
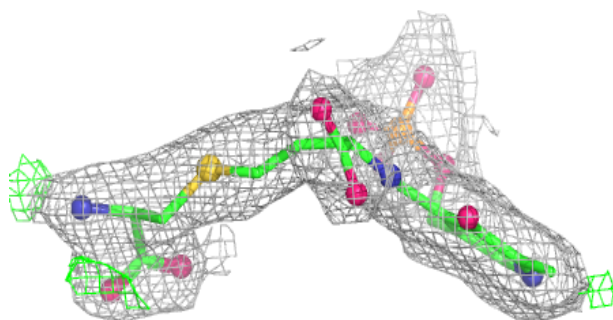
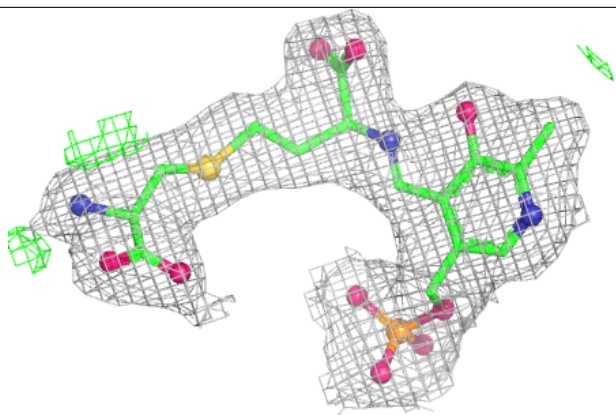
Electron density around E9U A 801:

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

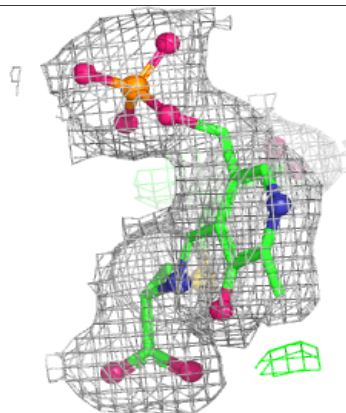
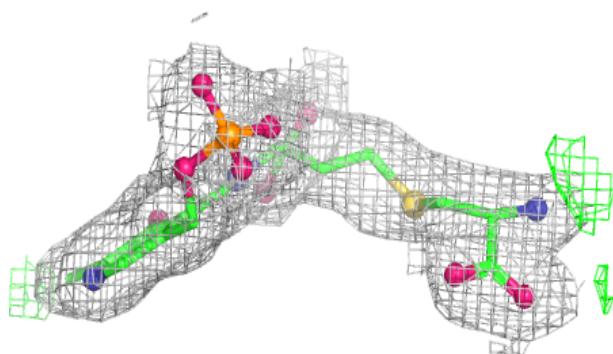
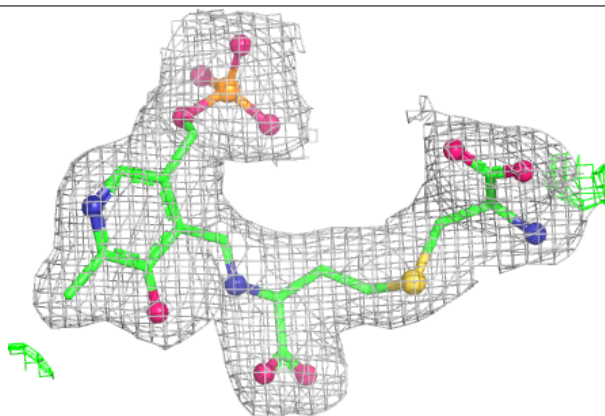


Electron density around E9U B 801:

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

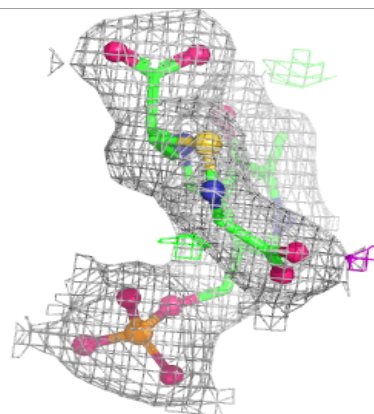
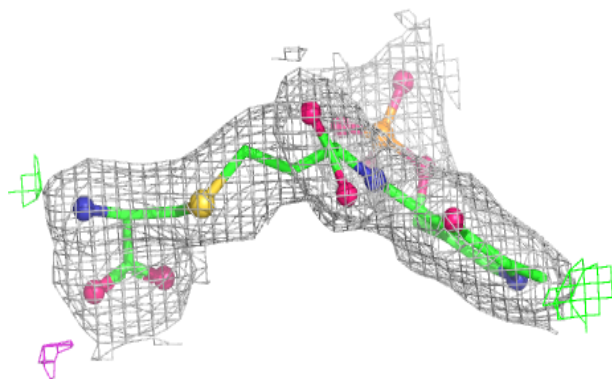
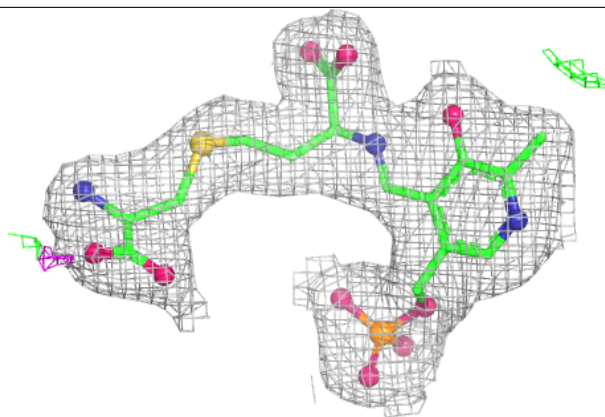
**Electron density around E9U C 801:**

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



Electron density around E9U D 801:

$2mF_o-DF_c$ (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.