



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 07:47 PM UTC

PDB ID : 4DPG / pdb_00004dpg
Title : Crystal Structure of Human LysRS: P38/AIMP2 Complex I
Authors : Fang, P.; Wang, J.; Bennett, S.P.; Guo, M.
Deposited on : 2012-02-13
Resolution : 2.84 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

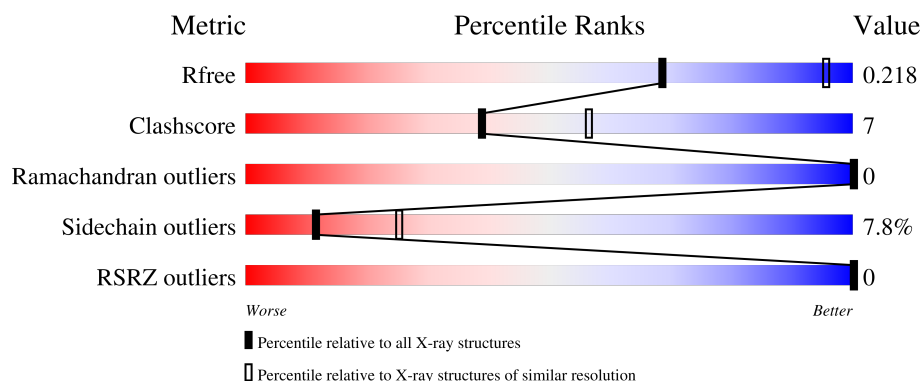
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.84 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	1520 (2.86-2.82)
Clashscore	190562	1559 (2.86-2.82)
Ramachandran outliers	187476	1517 (2.86-2.82)
Sidechain outliers	187428	1518 (2.86-2.82)
RSRZ outliers	180081	1521 (2.86-2.82)

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	513	
1	B	513	
1	C	513	
1	D	513	
1	E	513	

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Mol	Chain	Length	Quality of chain
1	F	513	80%17%..
1	G	513	76%19%..
1	H	513	77%18%..
2	I	54	37%19%.43%
2	J	54	39%15%.43%
2	K	54	39%15%.43%
2	L	54	35%20%.43%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 33740 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 Lysine-tRNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	501	Total	C	N	O	S	0	0	0
			3932	2519	668	718	27			
1	B	506	Total	C	N	O	S	0	0	0
			4055	2596	687	744	28			
1	C	501	Total	C	N	O	S	0	0	0
			3989	2553	677	731	28			
1	D	505	Total	C	N	O	S	0	0	0
			4031	2588	685	730	28			
1	E	500	Total	C	N	O	S	0	0	0
			3969	2540	671	730	28			
1	F	505	Total	C	N	O	S	0	0	0
			4023	2580	685	730	28			
1	G	501	Total	C	N	O	S	0	0	0
			3921	2507	667	719	28			
1	H	505	Total	C	N	O	S	0	0	0
			4045	2590	685	742	28			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	69	MET	-	expression tag	UNP Q15046
B	69	MET	-	expression tag	UNP Q15046
C	69	MET	-	expression tag	UNP Q15046
D	69	MET	-	expression tag	UNP Q15046
E	69	MET	-	expression tag	UNP Q15046
F	69	MET	-	expression tag	UNP Q15046
G	69	MET	-	expression tag	UNP Q15046
H	69	MET	-	expression tag	UNP Q15046

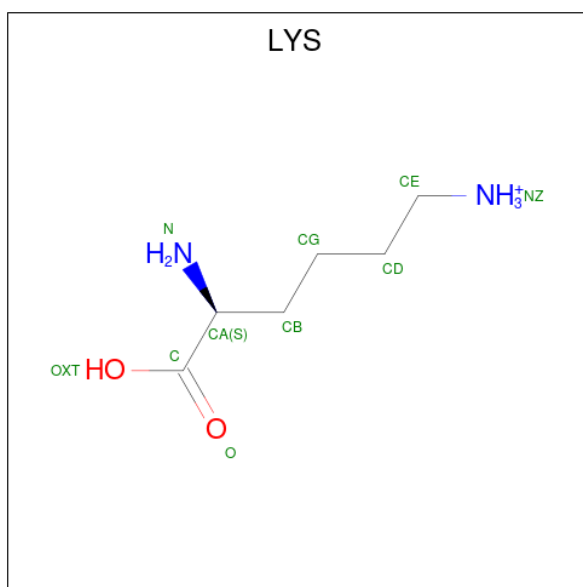
- Molecule 2 is a protein called Aminoacyl tRNA synthase complex-interacting multifunctional protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	I	31	Total	C	N	O	S	0	0	0
			224	144	40	37	3			
2	J	31	Total	C	N	O	S	0	0	0
			226	146	40	37	3			
2	K	31	Total	C	N	O	S	0	0	0
			224	144	40	37	3			
2	L	31	Total	C	N	O	S	0	0	0
			230	148	40	39	3			

There are 24 discrepancies between the modelled and reference sequences:

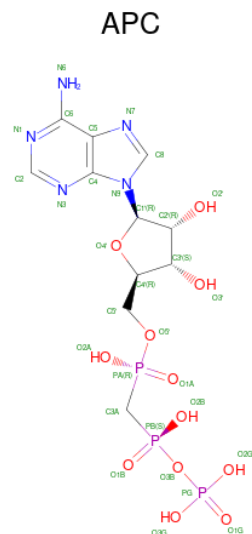
Chain	Residue	Modelled	Actual	Comment	Reference
I	49	HIS	-	expression tag	UNP Q13155
I	50	HIS	-	expression tag	UNP Q13155
I	51	HIS	-	expression tag	UNP Q13155
I	52	HIS	-	expression tag	UNP Q13155
I	53	HIS	-	expression tag	UNP Q13155
I	54	HIS	-	expression tag	UNP Q13155
J	49	HIS	-	expression tag	UNP Q13155
J	50	HIS	-	expression tag	UNP Q13155
J	51	HIS	-	expression tag	UNP Q13155
J	52	HIS	-	expression tag	UNP Q13155
J	53	HIS	-	expression tag	UNP Q13155
J	54	HIS	-	expression tag	UNP Q13155
K	49	HIS	-	expression tag	UNP Q13155
K	50	HIS	-	expression tag	UNP Q13155
K	51	HIS	-	expression tag	UNP Q13155
K	52	HIS	-	expression tag	UNP Q13155
K	53	HIS	-	expression tag	UNP Q13155
K	54	HIS	-	expression tag	UNP Q13155
L	49	HIS	-	expression tag	UNP Q13155
L	50	HIS	-	expression tag	UNP Q13155
L	51	HIS	-	expression tag	UNP Q13155
L	52	HIS	-	expression tag	UNP Q13155
L	53	HIS	-	expression tag	UNP Q13155
L	54	HIS	-	expression tag	UNP Q13155

- Molecule 3 is LYSINE (CCD ID: LYS) (formula: $C_6H_{15}N_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			10	6	2	2		
3	B	1	Total	C	N	O	0	0
			10	6	2	2		
3	C	1	Total	C	N	O	0	0
			10	6	2	2		
3	D	1	Total	C	N	O	0	0
			10	6	2	2		
3	E	1	Total	C	N	O	0	0
			10	6	2	2		
3	F	1	Total	C	N	O	0	0
			10	6	2	2		
3	G	1	Total	C	N	O	0	0
			10	6	2	2		
3	H	1	Total	C	N	O	0	0
			10	6	2	2		

- Molecule 4 is DIPHOSPHOMETHYLPHOSPHONIC ACID ADENOSYL ESTER (CCD ID: APC) (formula: $C_{11}H_{18}N_5O_{12}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 27	C 11	N 5	O 9	P 2	0	0
4	B	1	Total 31	C 11	N 5	O 12	P 3	0	0
4	C	1	Total 27	C 11	N 5	O 9	P 2	0	0
4	D	1	Total 31	C 11	N 5	O 12	P 3	0	0
4	E	1	Total 27	C 11	N 5	O 9	P 2	0	0
4	F	1	Total 31	C 11	N 5	O 12	P 3	0	0
4	G	1	Total 27	C 11	N 5	O 9	P 2	0	0
4	H	1	Total 31	C 11	N 5	O 12	P 3	0	0

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

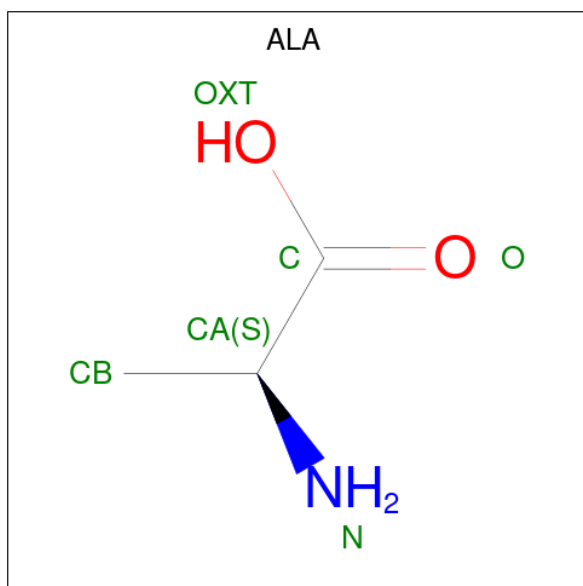
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Mg 1 1	0	0
5	B	1	Total Mg 1 1	0	0
5	C	1	Total Mg 1 1	0	0
5	D	1	Total Mg 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	E	1	Total	Mg	0	0
			1	1		
5	F	2	Total	Mg	0	0
			2	2		
5	G	1	Total	Mg	0	0
			1	1		
5	H	2	Total	Mg	0	0
			2	2		
5	J	1	Total	Mg	0	0
			1	1		
5	L	1	Total	Mg	0	0
			1	1		

- Molecule 6 is ALANINE (CCD ID: ALA) (formula: $C_3H_7NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	H	1	Total	C	N	O	0	0
			5	3	1	1		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	47	Total	O	0	0
			47	47		
7	B	85	Total	O	0	0
			85	85		

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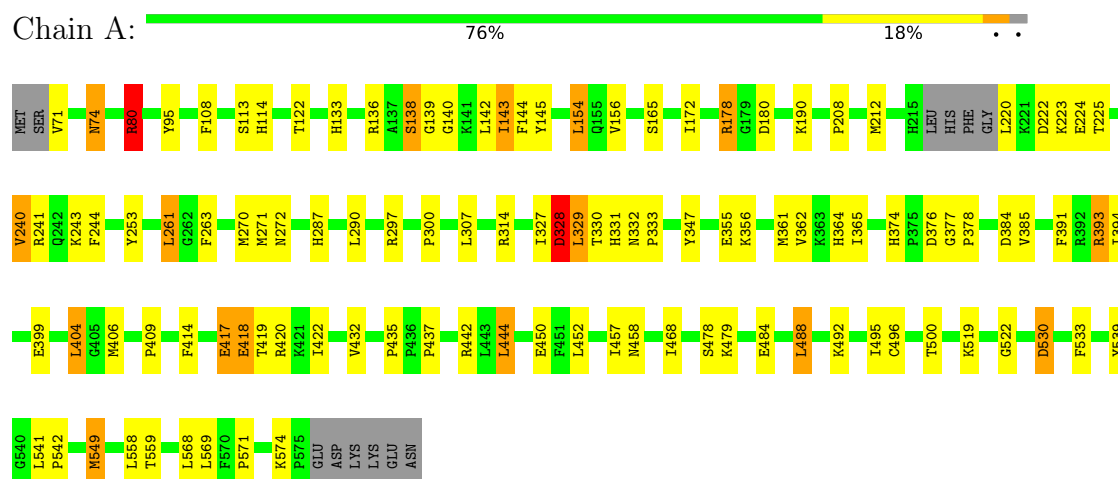
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	55	Total 55	O 55	0	0
7	D	71	Total 71	O 71	0	0
7	E	66	Total 66	O 66	0	0
7	F	81	Total 81	O 81	0	0
7	G	43	Total 43	O 43	0	0
7	H	62	Total 62	O 62	0	0
7	I	6	Total 6	O 6	0	0
7	J	3	Total 3	O 3	0	0
7	K	10	Total 10	O 10	0	0
7	L	13	Total 13	O 13	0	0

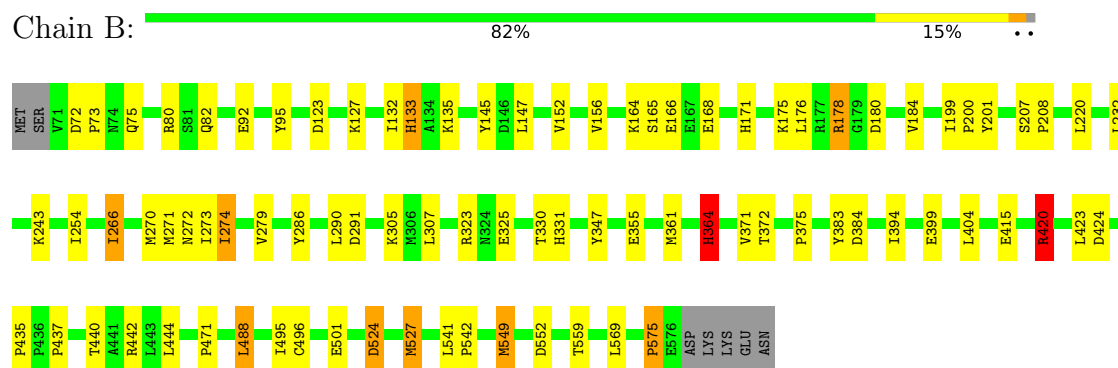
3 Residue-property plots

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.

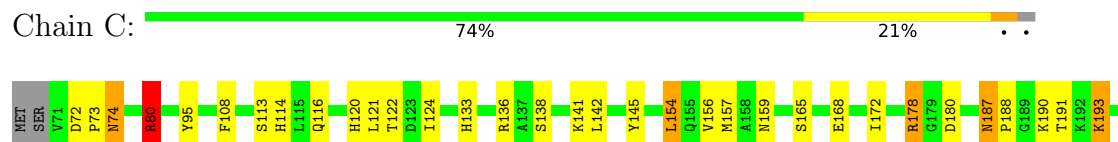
• Molecule 1: Lysine-tRNA ligase

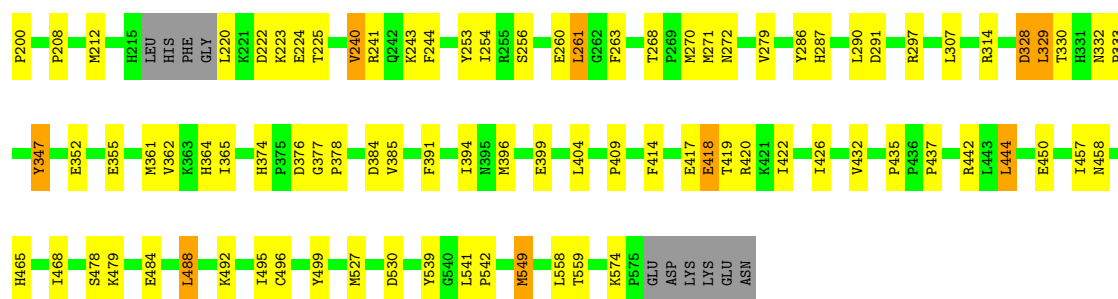


• Molecule 1: Lysine-tRNA ligase



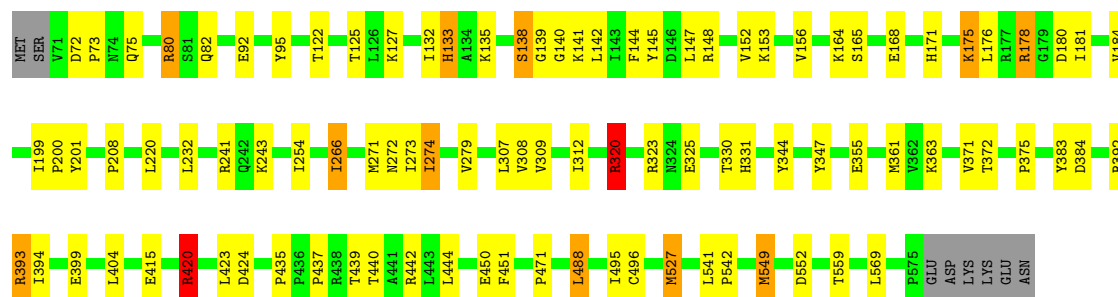
• Molecule 1: Lysine-tRNA ligase





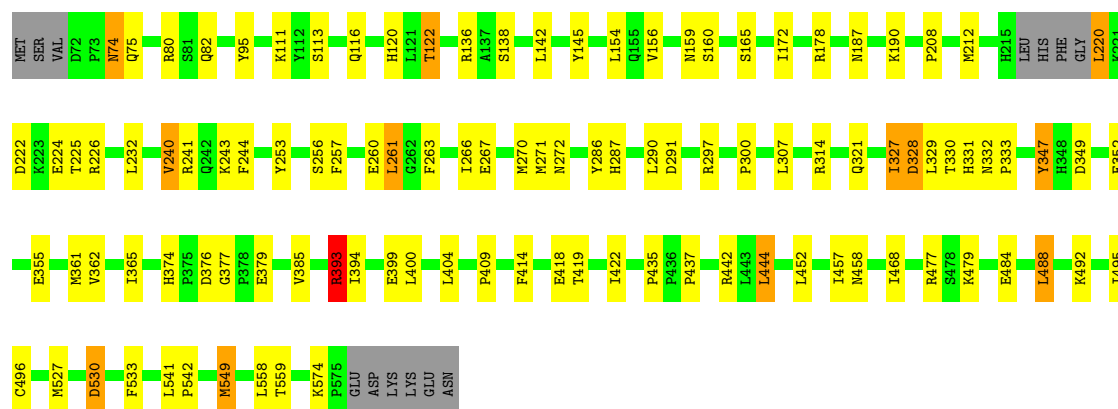
• Molecule 1: Lysine-tRNA ligase

Chain D: 80% 16%



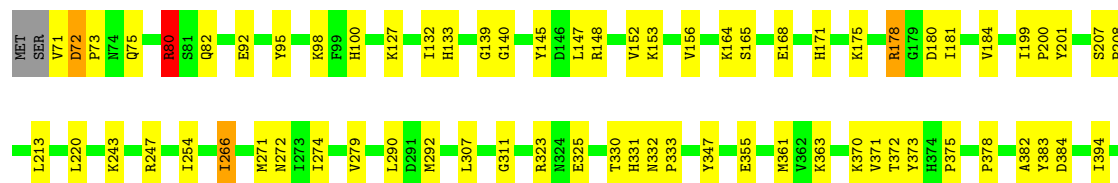
• Molecule 1: Lysine-tRNA ligase

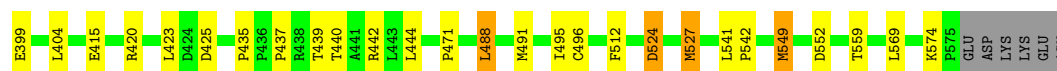
Chain E: 76% 19%



• Molecule 1: Lysine-tRNA ligase

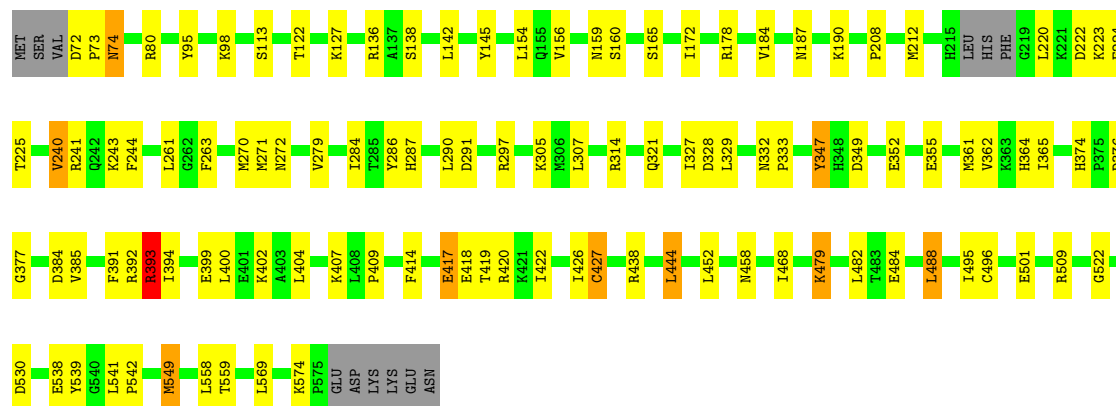
Chain F: 80% 17%





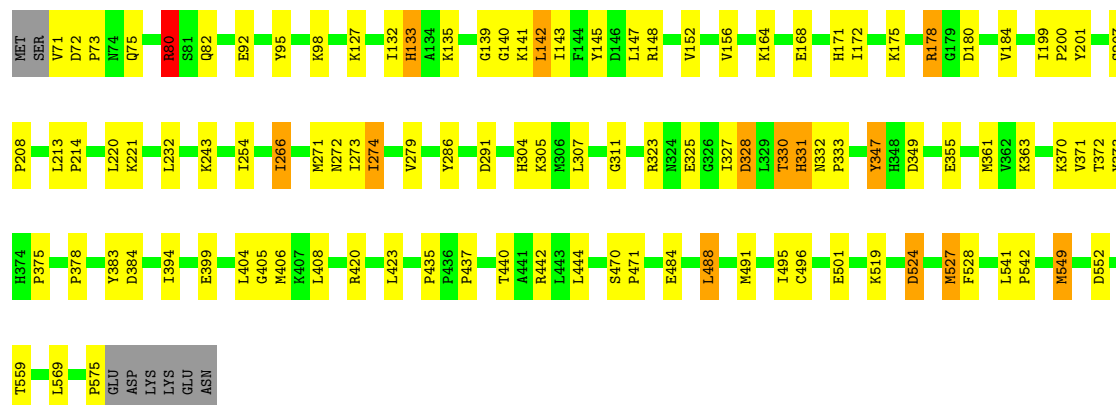
• Molecule 1: Lysine-tRNA ligase

Chain G: 76% 19%



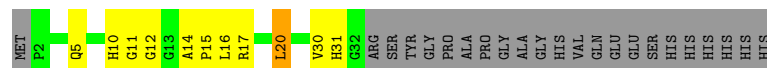
• Molecule 1: Lysine-tRNA ligase

Chain H: 77% 18%



• Molecule 2: Aminoacyl tRNA synthase complex-interacting multifunctional protein 2

Chain I: 37% 19% 43%

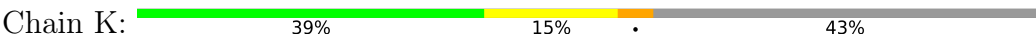


• Molecule 2: Aminoacyl tRNA synthase complex-interacting multifunctional protein 2

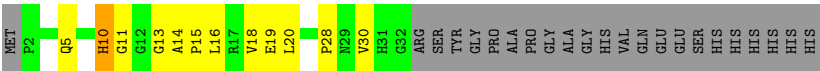
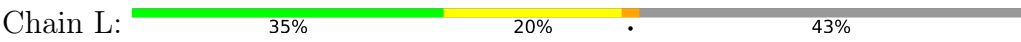
Chain J: 39% 15% 43%



- Molecule 2: Aminoacyl tRNA synthase complex-interacting multifunctional protein 2



- Molecule 2: Aminoacyl tRNA synthase complex-interacting multifunctional protein 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	79.20Å 122.01Å 149.20Å 89.16° 85.58° 89.71°	Depositor
Resolution (Å)	46.84 – 2.84 46.84 – 2.84	Depositor EDS
% Data completeness (in resolution range)	98.5 (46.84-2.84) 95.0 (46.84-2.84)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 2.86Å)	Xtriage
Refinement program	PHENIX 1.7.2_869	Depositor
R, R_{free}	0.185 , 0.223 0.180 , 0.218	Depositor DCC
R_{free} test set	6481 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.002	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 13.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.339 for -h,k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	33740	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% 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: MG, APC

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	0.64	1/4025 (0.0%)	0.95	10/5457 (0.2%)
1	B	0.63	0/4152	0.94	13/5619 (0.2%)
1	C	0.61	0/4082	0.90	6/5526 (0.1%)
1	D	0.65	0/4128	1.00	16/5586 (0.3%)
1	E	0.61	0/4062	0.92	12/5503 (0.2%)
1	F	0.63	0/4120	0.92	10/5577 (0.2%)
1	G	0.60	0/4012	0.92	8/5440 (0.1%)
1	H	0.66	1/4142 (0.0%)	0.93	8/5607 (0.1%)
2	I	0.64	0/233	1.11	1/319 (0.3%)
2	J	0.81	1/235 (0.4%)	1.13	2/322 (0.6%)
2	K	0.83	0/233	1.19	2/319 (0.6%)
2	L	0.84	1/239 (0.4%)	1.07	0/327
All	All	0.63	4/33663 (0.0%)	0.94	88/45602 (0.2%)

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	A	0	1
1	B	0	1
1	C	0	1
1	E	0	1
1	G	0	1
All	All	0	5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	241	ARG	CD-NE	-5.20	1.39	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	331	HIS	CG-ND1	-5.18	1.32	1.38
2	L	10	HIS	CA-C	-5.08	1.47	1.52
2	J	31	HIS	CG-ND1	-5.04	1.32	1.38

All (88) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	320	ARG	NE-CZ-NH2	13.34	131.20	119.20
1	A	241	ARG	NE-CZ-NH2	12.68	130.61	119.20
1	D	393	ARG	NE-CZ-NH2	12.04	130.03	119.20
1	D	320	ARG	NE-CZ-NH1	-10.70	110.80	121.50
1	A	241	ARG	NE-CZ-NH1	-10.58	110.92	121.50
1	D	178	ARG	NE-CZ-NH2	10.17	128.35	119.20
1	D	393	ARG	NE-CZ-NH1	-9.80	111.70	121.50
1	B	178	ARG	NE-CZ-NH2	9.69	127.92	119.20
1	D	320	ARG	CD-NE-CZ	9.62	137.88	124.40
1	E	178	ARG	NE-CZ-NH2	9.38	127.64	119.20
1	H	420	ARG	NE-CZ-NH2	9.20	127.48	119.20
1	G	178	ARG	NE-CZ-NH2	9.18	127.46	119.20
1	G	393	ARG	NE-CZ-NH2	9.13	127.41	119.20
1	F	420	ARG	NE-CZ-NH2	8.78	127.11	119.20
2	I	17	ARG	N-CA-C	-8.71	102.56	113.02
1	A	241	ARG	CD-NE-CZ	8.65	136.50	124.40
1	E	477	ARG	NE-CZ-NH2	-8.64	111.42	119.20
1	E	393	ARG	NE-CZ-NH2	8.53	126.88	119.20
1	C	80	ARG	NE-CZ-NH2	8.25	126.63	119.20
1	E	477	ARG	NE-CZ-NH1	8.22	129.72	121.50
1	A	80	ARG	NE-CZ-NH2	7.97	126.37	119.20
1	D	393	ARG	CD-NE-CZ	7.83	135.36	124.40
1	E	477	ARG	CD-NE-CZ	7.75	135.24	124.40
1	F	420	ARG	NE-CZ-NH1	-7.46	114.04	121.50
1	H	420	ARG	NE-CZ-NH1	-7.18	114.32	121.50
1	F	80	ARG	NE-CZ-NH2	7.10	125.59	119.20
1	B	178	ARG	NE-CZ-NH1	-7.07	114.43	121.50
1	H	80	ARG	NE-CZ-NH2	7.02	125.52	119.20
1	E	178	ARG	NE-CZ-NH1	-6.98	114.52	121.50
1	D	178	ARG	NE-CZ-NH1	-6.91	114.59	121.50
1	H	331	HIS	N-CA-C	6.86	120.35	108.76
1	G	178	ARG	NE-CZ-NH1	-6.77	114.73	121.50
1	G	393	ARG	NE-CZ-NH1	-6.69	114.81	121.50
1	A	80	ARG	NE-CZ-NH1	-6.60	114.90	121.50
1	D	138	SER	N-CA-C	-6.58	103.74	113.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	420	ARG	CD-NE-CZ	6.57	133.59	124.40
2	J	16	LEU	N-CA-C	-6.53	103.81	113.40
2	J	17	ARG	N-CA-C	-6.36	105.39	112.57
1	E	393	ARG	NE-CZ-NH1	-6.32	115.18	121.50
1	F	420	ARG	CD-NE-CZ	6.20	133.07	124.40
1	B	364	HIS	CG-CD2-NE2	6.18	113.39	107.20
1	D	420	ARG	NE-CZ-NH1	6.15	127.65	121.50
1	D	527	MET	N-CA-C	6.11	119.51	110.46
1	F	178	ARG	CD-NE-CZ	6.08	132.91	124.40
1	H	178	ARG	CD-NE-CZ	6.05	132.88	124.40
1	C	80	ARG	NE-CZ-NH1	-6.04	115.46	121.50
1	D	420	ARG	NE-CZ-NH2	-6.03	113.77	119.20
1	E	477	ARG	CB-CG-CD	-5.95	97.62	111.30
1	D	178	ARG	CD-NE-CZ	5.89	132.65	124.40
1	G	178	ARG	CD-NE-CZ	5.87	132.61	124.40
1	D	420	ARG	CD-NE-CZ	5.86	132.60	124.40
1	B	178	ARG	CD-NE-CZ	5.85	132.59	124.40
1	B	420	ARG	NE-CZ-NH1	5.85	127.35	121.50
1	B	420	ARG	CD-NE-CZ	5.84	132.57	124.40
1	H	527	MET	N-CA-C	5.81	119.06	110.46
1	E	178	ARG	CD-NE-CZ	5.80	132.51	124.40
1	F	527	MET	N-CA-C	5.78	119.02	110.46
1	B	331	HIS	N-CA-C	5.70	117.71	108.26
1	B	364	HIS	ND1-CG-CD2	-5.69	100.41	106.10
1	B	527	MET	N-CA-C	5.67	118.85	110.46
1	A	178	ARG	NE-CZ-NH1	5.57	127.07	121.50
1	C	178	ARG	NE-CZ-NH1	5.54	127.04	121.50
1	F	331	HIS	N-CA-C	5.50	117.40	108.26
1	A	328	ASP	N-CA-C	5.46	115.77	108.23
1	C	499	TYR	N-CA-C	5.45	117.13	108.79
1	D	331	HIS	N-CA-C	5.45	117.31	108.26
1	A	178	ARG	CD-NE-CZ	5.41	131.97	124.40
1	B	575	PRO	N-CA-C	5.40	119.05	111.33
1	A	522	GLY	N-CA-C	5.35	124.14	114.46
2	K	11	GLY	N-CA-C	-5.34	103.16	111.12
1	B	420	ARG	NE-CZ-NH2	-5.32	114.41	119.20
1	C	178	ARG	CD-NE-CZ	5.30	131.82	124.40
1	E	393	ARG	CD-NE-CZ	5.30	131.82	124.40
1	D	176	LEU	N-CA-C	5.27	117.87	109.96
1	F	80	ARG	NE-CZ-NH1	-5.21	116.29	121.50
1	C	527	MET	N-CA-C	5.13	117.48	110.55
1	B	176	LEU	N-CA-C	5.10	117.61	109.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	527	MET	N-CA-C	5.05	117.20	110.53
1	G	393	ARG	CD-NE-CZ	5.05	131.47	124.40
1	A	393	ARG	CD-NE-CZ	5.04	131.46	124.40
1	B	166	GLU	N-CA-C	-5.04	105.86	111.36
1	E	160	SER	N-CA-C	5.04	116.77	111.28
1	F	72	ASP	CA-C-N	5.02	124.80	119.28
1	F	72	ASP	C-N-CA	5.02	124.80	119.28
1	H	80	ARG	NE-CZ-NH1	-5.01	116.48	121.50
1	G	522	GLY	N-CA-C	5.01	123.53	114.46
1	G	160	SER	N-CA-C	5.00	116.74	111.28
2	K	16	LEU	N-CA-C	-5.00	103.11	113.31

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	574	LYS	Peptide
1	B	575	PRO	Peptide
1	C	574	LYS	Peptide
1	E	574	LYS	Peptide
1	G	574	LYS	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	3932	0	3804	60	0
1	B	4055	0	3984	42	0
1	C	3989	0	3910	75	0
1	D	4031	0	3977	51	0
1	E	3969	0	3873	63	0
1	F	4023	0	3964	47	0
1	G	3921	0	3801	51	0
1	H	4045	0	3981	63	0
2	I	224	0	200	6	0
2	J	226	0	207	3	0
2	K	224	0	200	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	L	230	0	211	15	0
3	A	10	0	12	0	0
3	B	10	0	12	0	0
3	C	10	0	12	0	0
3	D	10	0	12	0	0
3	E	10	0	12	0	0
3	F	10	0	12	0	0
3	G	10	0	12	0	0
3	H	10	0	12	0	0
4	A	27	0	14	1	0
4	B	31	0	14	0	0
4	C	27	0	14	1	0
4	D	31	0	14	0	0
4	E	27	0	14	2	0
4	F	31	0	14	0	0
4	G	27	0	14	0	0
4	H	31	0	14	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0
5	F	2	0	0	0	0
5	G	1	0	0	0	0
5	H	2	0	0	0	0
5	J	1	0	0	0	0
5	L	1	0	0	0	0
6	H	5	0	4	1	0
7	A	47	0	0	2	0
7	B	85	0	0	3	0
7	C	55	0	0	4	0
7	D	71	0	0	3	0
7	E	66	0	0	1	0
7	F	81	0	0	2	0
7	G	43	0	0	1	0
7	H	62	0	0	3	0
7	I	6	0	0	0	0
7	J	3	0	0	0	0
7	K	10	0	0	0	0
7	L	13	0	0	1	0
All	All	33740	0	32324	446	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 7.

All (446) 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:270:MET:HE1	7:C:739:HOH:O	1.38	1.16
1:C:124:ILE:O	1:C:187:ASN:HB3	1.68	0.92
7:B:776:HOH:O	2:I:14:ALA:HB2	1.72	0.88
1:C:256:SER:OG	2:K:14:ALA:HB1	1.73	0.88
1:G:427:CYS:SG	1:G:438:ARG:NH1	2.51	0.84
1:B:394:ILE:HB	1:B:399:GLU:HG3	1.64	0.80
1:E:120:HIS:CD2	1:E:122:THR:HG23	2.17	0.78
1:H:394:ILE:HB	1:H:399:GLU:HG3	1.65	0.77
1:F:394:ILE:HB	1:F:399:GLU:HG3	1.66	0.77
1:E:287:HIS:CD2	1:E:290:LEU:HG	2.21	0.76
1:D:394:ILE:HB	1:D:399:GLU:HG3	1.66	0.75
1:A:327:ILE:HD12	1:B:290:LEU:HD11	1.68	0.75
1:C:287:HIS:HD2	1:C:290:LEU:HG	1.49	0.75
1:C:287:HIS:CD2	1:C:290:LEU:HG	2.22	0.75
1:E:328:ASP:OD1	1:E:328:ASP:C	2.30	0.73
1:E:287:HIS:HD2	1:E:290:LEU:HG	1.51	0.73
1:G:409:PRO:HG2	1:G:419:THR:HG23	1.69	0.73
1:G:287:HIS:CD2	1:G:290:LEU:HG	2.24	0.73
2:J:15:PRO:O	2:J:16:LEU:CB	2.35	0.73
1:A:409:PRO:HG2	1:A:419:THR:HG23	1.70	0.72
2:I:15:PRO:O	2:I:16:LEU:CB	2.37	0.71
1:C:409:PRO:HG2	1:C:419:THR:HG23	1.71	0.71
1:G:287:HIS:HD2	1:G:290:LEU:HG	1.56	0.71
1:B:435:PRO:O	1:B:442:ARG:NH2	2.24	0.70
1:H:575:PRO:C	6:H:601:ALA:N	2.50	0.70
1:B:364:HIS:C	1:B:364:HIS:HD1	2.01	0.69
1:C:133:HIS:O	1:C:178:ARG:HD2	1.92	0.69
1:C:193:LYS:HE3	1:C:193:LYS:H	1.58	0.69
1:D:435:PRO:O	1:D:442:ARG:NH2	2.27	0.68
1:H:488:LEU:HB2	1:H:496:CYS:HB2	1.75	0.68
1:C:256:SER:HB3	2:K:15:PRO:O	1.94	0.68
1:E:253:TYR:HD1	2:L:15:PRO:HB3	1.58	0.68
1:E:409:PRO:HG2	1:E:419:THR:HG23	1.75	0.67
1:B:133:HIS:O	1:B:178:ARG:HD2	1.94	0.67
1:A:133:HIS:O	1:A:178:ARG:HD2	1.94	0.66
1:G:113:SER:O	1:G:190:LYS:NZ	2.27	0.66
1:D:241:ARG:NH2	7:D:727:HOH:O	2.29	0.66
1:B:488:LEU:HB2	1:B:496:CYS:HB2	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:253:TYR:CD1	2:L:15:PRO:HB3	2.32	0.65
1:E:328:ASP:OD1	1:E:330:THR:N	2.29	0.65
1:G:328:ASP:OD1	1:G:329:LEU:N	2.30	0.65
2:K:15:PRO:O	2:K:16:LEU:CB	2.42	0.65
2:L:11:GLY:HA2	7:L:211:HOH:O	1.96	0.65
1:F:435:PRO:O	1:F:442:ARG:NH2	2.30	0.65
1:E:328:ASP:OD1	1:E:329:LEU:N	2.30	0.64
1:C:328:ASP:OD1	1:C:329:LEU:N	2.29	0.64
1:E:361:MET:HE1	1:E:549:MET:HE1	1.78	0.64
1:E:120:HIS:HD2	1:E:122:THR:HG23	1.62	0.64
1:H:133:HIS:HB3	7:H:739:HOH:O	1.96	0.64
1:F:488:LEU:HB2	1:F:496:CYS:HB2	1.77	0.64
1:A:328:ASP:OD1	1:A:329:LEU:N	2.30	0.64
1:E:113:SER:O	1:E:190:LYS:NZ	2.31	0.64
1:H:328:ASP:OD1	1:H:331:HIS:ND1	2.30	0.64
1:C:116:GLN:HE22	1:F:382:ALA:H	1.43	0.64
1:C:74:ASN:N	1:C:74:ASN:OD1	2.31	0.63
1:C:253:TYR:HD1	2:K:15:PRO:HB3	1.63	0.63
1:G:541:LEU:HD12	1:G:542:PRO:HD2	1.80	0.63
1:E:256:SER:OG	2:L:15:PRO:HD2	1.98	0.63
1:D:541:LEU:HD12	1:D:542:PRO:HD2	1.79	0.63
1:H:435:PRO:O	1:H:442:ARG:NH2	2.32	0.63
1:A:253:TYR:HD1	1:A:364:HIS:HD2	1.47	0.63
1:A:541:LEU:HD12	1:A:542:PRO:HD2	1.81	0.62
1:B:541:LEU:HD12	1:B:542:PRO:HD2	1.81	0.62
1:A:74:ASN:OD1	1:A:74:ASN:N	2.31	0.62
1:E:541:LEU:HD12	1:E:542:PRO:HD2	1.80	0.62
1:G:74:ASN:OD1	1:G:74:ASN:N	2.30	0.62
1:A:113:SER:O	1:A:190:LYS:NZ	2.33	0.62
1:B:164:LYS:HB3	1:B:201:TYR:CZ	2.35	0.62
1:B:273:ILE:HG13	1:B:274:ILE:HG12	1.80	0.62
1:C:488:LEU:HB2	1:C:496:CYS:HB2	1.81	0.62
1:A:114:HIS:NE2	1:H:378:PRO:O	2.31	0.62
1:A:328:ASP:OD1	1:A:330:THR:N	2.30	0.62
1:A:328:ASP:OD2	1:A:331:HIS:ND1	2.30	0.62
1:D:139:GLY:HA3	1:D:140:GLY:C	2.24	0.62
1:H:273:ILE:HG13	1:H:274:ILE:HG12	1.82	0.62
1:H:328:ASP:OD1	1:H:331:HIS:CG	2.53	0.62
1:D:133:HIS:O	1:D:178:ARG:HD2	1.98	0.61
1:G:361:MET:HE1	1:G:549:MET:HE1	1.82	0.61
1:C:253:TYR:CD1	2:K:15:PRO:HB3	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:74:ASN:OD1	1:E:74:ASN:N	2.33	0.61
1:D:273:ILE:HG13	1:D:274:ILE:HG12	1.81	0.60
1:C:361:MET:HE1	1:C:549:MET:HE1	1.83	0.60
1:H:524:ASP:OD2	1:H:524:ASP:N	2.28	0.60
1:C:80:ARG:NH1	1:C:180:ASP:OD2	2.33	0.60
1:C:113:SER:O	1:C:190:LYS:NZ	2.34	0.60
1:F:95:TYR:CE2	1:F:208:PRO:HD2	2.37	0.59
1:G:488:LEU:HB2	1:G:496:CYS:HB2	1.83	0.59
1:F:80:ARG:NH1	1:F:180:ASP:OD1	2.36	0.59
1:F:541:LEU:HD12	1:F:542:PRO:HD2	1.83	0.59
1:A:80:ARG:NH1	1:A:180:ASP:OD2	2.34	0.59
1:A:488:LEU:HB2	1:A:496:CYS:HB2	1.84	0.59
1:D:139:GLY:CA	1:D:140:GLY:C	2.75	0.59
1:H:95:TYR:CE2	1:H:208:PRO:HD2	2.37	0.59
1:C:328:ASP:OD1	1:C:330:THR:N	2.30	0.59
1:H:541:LEU:HD12	1:H:542:PRO:HD2	1.84	0.59
1:C:253:TYR:HD1	1:C:364:HIS:HD2	1.51	0.59
1:A:270:MET:HE1	7:B:770:HOH:O	2.01	0.58
1:D:95:TYR:CE2	1:D:208:PRO:HD2	2.38	0.58
1:A:488:LEU:HD13	1:A:495:ILE:HD11	1.86	0.58
1:E:327:ILE:O	1:E:327:ILE:HG22	2.02	0.58
1:E:488:LEU:HD13	1:E:495:ILE:HD11	1.86	0.58
1:E:488:LEU:HB2	1:E:496:CYS:HB2	1.86	0.58
1:G:145:TYR:HB2	1:G:156:VAL:HB	1.86	0.58
1:B:364:HIS:C	1:B:364:HIS:ND1	2.62	0.58
1:B:286:TYR:OH	1:B:291:ASP:OD1	2.21	0.57
1:E:270:MET:HA	1:E:297:ARG:HD3	1.86	0.57
1:D:488:LEU:HB2	1:D:496:CYS:HB2	1.85	0.57
7:E:762:HOH:O	2:L:19:GLU:HG3	2.05	0.57
1:G:270:MET:HA	1:G:297:ARG:HD3	1.87	0.57
1:H:164:LYS:HB3	1:H:201:TYR:CZ	2.38	0.57
1:C:541:LEU:HD12	1:C:542:PRO:HD2	1.86	0.57
1:C:145:TYR:HB2	1:C:156:VAL:HB	1.87	0.57
1:C:142:LEU:HD13	1:C:159:ASN:HB3	1.87	0.57
1:F:323:ARG:NH1	1:F:325:GLU:OE2	2.37	0.57
1:B:95:TYR:CE2	1:B:208:PRO:HD2	2.40	0.56
1:A:361:MET:HE1	1:A:549:MET:HE1	1.87	0.56
1:B:323:ARG:NH1	1:B:325:GLU:OE2	2.37	0.56
1:H:323:ARG:NH1	1:H:325:GLU:OE2	2.36	0.56
1:E:256:SER:OG	2:L:14:ALA:HB1	2.06	0.56
1:A:145:TYR:HB2	1:A:156:VAL:HB	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:145:TYR:HB2	1:E:156:VAL:HB	1.87	0.56
1:G:142:LEU:HD13	1:G:159:ASN:HB3	1.87	0.56
1:A:253:TYR:CD1	1:A:364:HIS:HD2	2.23	0.56
1:F:164:LYS:HB3	1:F:201:TYR:CZ	2.41	0.56
1:D:323:ARG:NH1	1:D:325:GLU:OE2	2.38	0.56
1:A:138:SER:HB2	1:A:142:LEU:HD23	1.88	0.55
1:E:142:LEU:HD13	1:E:159:ASN:HB3	1.88	0.55
1:H:254:ILE:HD11	1:H:549:MET:CE	2.36	0.55
1:B:270:MET:HE1	7:B:770:HOH:O	2.06	0.55
1:H:286:TYR:OH	1:H:291:ASP:OD1	2.20	0.55
1:F:361:MET:HE1	1:F:549:MET:HE1	1.89	0.54
1:E:256:SER:HB3	2:L:15:PRO:O	2.07	0.54
1:F:145:TYR:HB2	1:F:156:VAL:HB	1.89	0.54
1:H:80:ARG:NH1	1:H:180:ASP:OD1	2.40	0.54
1:G:287:HIS:HD2	1:G:290:LEU:CG	2.21	0.54
1:C:120:HIS:NE2	1:C:187:ASN:OD1	2.40	0.54
1:H:488:LEU:HD13	1:H:495:ILE:HD11	1.89	0.54
1:D:361:MET:HE1	1:D:549:MET:HE1	1.90	0.53
1:H:423:LEU:HD12	1:H:440:THR:HG23	1.90	0.53
1:F:471:PRO:HG2	1:F:527:MET:HE3	1.90	0.53
1:E:307:LEU:HD11	1:F:569:LEU:HD13	1.91	0.53
1:C:488:LEU:HD13	1:C:495:ILE:HD11	1.90	0.53
1:C:240:VAL:HG12	1:C:244:PHE:HE2	1.74	0.53
1:C:364:HIS:CD2	2:K:15:PRO:HB3	2.44	0.53
1:D:164:LYS:HB3	1:D:201:TYR:CZ	2.43	0.53
1:E:240:VAL:HG12	1:E:244:PHE:HE2	1.73	0.53
1:H:145:TYR:HB2	1:H:156:VAL:HB	1.90	0.53
1:C:253:TYR:CD1	1:C:364:HIS:HD2	2.27	0.53
1:F:132:ILE:HD12	1:F:180:ASP:HB2	1.91	0.53
1:C:116:GLN:NE2	1:F:382:ALA:H	2.07	0.52
1:C:270:MET:HA	1:C:297:ARG:HD3	1.90	0.52
1:D:471:PRO:HG2	1:D:527:MET:HE3	1.91	0.52
1:A:222:ASP:OD1	1:A:225:THR:OG1	2.27	0.52
1:G:417:GLU:OE2	1:G:420:ARG:NE	2.42	0.52
1:A:332:ASN:CG	1:A:333:PRO:HD2	2.35	0.52
1:A:270:MET:HA	1:A:297:ARG:HD3	1.91	0.51
1:D:92:GLU:OE1	1:D:171:HIS:NE2	2.42	0.51
1:E:435:PRO:O	1:E:442:ARG:NH2	2.43	0.51
1:H:142:LEU:HD12	1:H:143:ILE:N	2.25	0.51
1:G:222:ASP:OD1	1:G:225:THR:OG1	2.28	0.51
1:G:427:CYS:HG	1:G:438:ARG:NH1	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:ILE:HG22	1:A:144:PHE:N	2.25	0.51
1:A:243:LYS:HG2	1:A:559:THR:HB	1.92	0.51
1:A:394:ILE:HD12	1:A:452:LEU:HD22	1.93	0.51
1:B:92:GLU:OE1	1:B:171:HIS:NE2	2.43	0.51
1:B:145:TYR:HB2	1:B:156:VAL:HB	1.91	0.51
1:A:240:VAL:HG12	1:A:244:PHE:HE2	1.76	0.51
1:A:287:HIS:CE1	1:A:290:LEU:HG	2.45	0.51
1:B:254:ILE:HD11	1:B:549:MET:CE	2.41	0.51
1:B:471:PRO:HG2	1:B:527:MET:HE3	1.93	0.51
1:D:423:LEU:HD12	1:D:440:THR:HG23	1.92	0.51
1:A:307:LEU:HD11	1:B:569:LEU:HD13	1.93	0.51
1:B:132:ILE:HD12	1:B:180:ASP:HB2	1.92	0.50
1:C:243:LYS:HG2	1:C:559:THR:HB	1.93	0.50
1:E:253:TYR:CD1	2:L:15:PRO:CB	2.94	0.50
1:G:307:LEU:HD11	1:H:569:LEU:HD13	1.94	0.50
1:G:422:ILE:O	1:G:426:ILE:HG13	2.11	0.50
1:G:488:LEU:HD13	1:G:495:ILE:HD11	1.92	0.50
1:B:123:ASP:HB3	1:F:425:ASP:HB3	1.93	0.50
1:D:145:TYR:HB2	1:D:156:VAL:HB	1.94	0.50
1:A:287:HIS:ND1	1:A:290:LEU:HG	2.27	0.50
1:A:327:ILE:CD1	1:B:290:LEU:HD11	2.39	0.50
1:E:347:TYR:CE1	1:E:484:GLU:HB3	2.46	0.50
1:F:254:ILE:HD11	1:F:549:MET:CE	2.41	0.50
1:G:364:HIS:C	1:G:364:HIS:HD1	2.20	0.50
1:H:471:PRO:HG2	1:H:527:MET:HE3	1.94	0.50
1:B:423:LEU:HD12	1:B:440:THR:HG23	1.94	0.50
1:E:95:TYR:CE2	1:E:208:PRO:HD2	2.46	0.50
1:G:240:VAL:HG12	1:G:244:PHE:HE2	1.76	0.50
1:H:92:GLU:OE1	1:H:171:HIS:NE2	2.45	0.50
1:A:417:GLU:OE2	1:A:420:ARG:NE	2.43	0.50
1:E:222:ASP:OD1	1:E:225:THR:OG1	2.27	0.50
1:C:95:TYR:CE2	1:C:208:PRO:HD2	2.47	0.50
1:G:347:TYR:CE1	1:G:484:GLU:HB3	2.47	0.50
1:E:330:THR:O	1:E:330:THR:OG1	2.30	0.49
1:E:385:VAL:HG13	1:E:458:ASN:HA	1.94	0.49
1:H:132:ILE:HD12	1:H:180:ASP:HB2	1.93	0.49
1:F:92:GLU:OE1	1:F:171:HIS:NE2	2.45	0.49
1:F:488:LEU:HD13	1:F:495:ILE:HD11	1.93	0.49
1:H:361:MET:HE1	1:H:549:MET:HE1	1.94	0.49
1:A:457:ILE:HD11	1:A:492:LYS:HE3	1.95	0.49
1:E:253:TYR:HA	2:L:15:PRO:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:95:TYR:CE2	1:G:208:PRO:HD2	2.47	0.49
1:C:261:LEU:HD13	2:K:20:LEU:HD22	1.95	0.49
1:C:417:GLU:OE2	1:C:420:ARG:NE	2.43	0.49
1:D:175:LYS:HE3	7:D:723:HOH:O	2.12	0.49
1:H:98:LYS:HA	7:H:706:HOH:O	2.13	0.49
1:B:524:ASP:OD2	1:B:524:ASP:N	2.26	0.49
1:C:362:VAL:HG22	1:C:558:LEU:HD21	1.93	0.49
1:D:254:ILE:HD11	1:D:549:MET:CE	2.43	0.49
1:C:268:THR:O	1:D:320:ARG:NH1	2.36	0.48
1:H:330:THR:O	1:H:331:HIS:HD2	1.96	0.48
1:A:435:PRO:O	1:A:442:ARG:NH2	2.46	0.48
1:B:437:PRO:O	1:B:442:ARG:HD3	2.12	0.48
1:G:243:LYS:HG2	1:G:559:THR:HB	1.95	0.48
1:C:364:HIS:HD2	2:K:15:PRO:HB3	1.78	0.48
1:C:418:GLU:O	1:C:422:ILE:HG13	2.14	0.48
1:D:437:PRO:O	1:D:442:ARG:HD3	2.14	0.48
1:B:305:LYS:NZ	1:B:501:GLU:OE1	2.38	0.48
1:C:114:HIS:NE2	1:F:378:PRO:O	2.44	0.48
1:C:222:ASP:OD1	1:C:225:THR:OG1	2.30	0.48
1:G:394:ILE:HD12	1:G:452:LEU:HD22	1.96	0.48
1:A:138:SER:CB	1:A:142:LEU:HD23	2.44	0.48
7:A:726:HOH:O	2:I:31:HIS:HE1	1.97	0.48
1:C:253:TYR:CD1	2:K:15:PRO:CB	2.96	0.48
1:G:385:VAL:HG13	1:G:458:ASN:HA	1.96	0.48
1:H:140:GLY:O	1:H:141:LYS:CB	2.60	0.48
1:C:435:PRO:O	1:C:442:ARG:NH2	2.46	0.48
1:E:418:GLU:O	1:E:422:ILE:HG13	2.14	0.48
1:E:120:HIS:CD2	1:E:122:THR:CG2	2.95	0.47
1:A:385:VAL:HG13	1:A:458:ASN:HA	1.95	0.47
1:D:148:ARG:HD3	1:D:153:LYS:HB3	1.96	0.47
1:G:355:GLU:OE2	1:G:393:ARG:NH2	2.47	0.47
1:A:488:LEU:CB	1:A:496:CYS:HB2	2.43	0.47
1:G:241:ARG:NH2	1:H:311:GLY:O	2.48	0.47
1:A:347:TYR:CE1	1:A:484:GLU:HB3	2.49	0.47
1:E:394:ILE:HB	1:E:399:GLU:HG3	1.96	0.47
1:F:423:LEU:HD12	1:F:440:THR:HG23	1.96	0.47
1:H:328:ASP:OD1	1:H:331:HIS:HB2	2.15	0.47
1:A:418:GLU:O	1:A:422:ILE:HG13	2.15	0.47
1:D:140:GLY:O	1:D:141:LYS:HB2	2.15	0.47
1:E:331:HIS:HD2	4:E:602:APC:C6	2.27	0.47
1:G:263:PHE:CD2	1:G:314:ARG:HB3	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:444:LEU:HA	1:D:444:LEU:HD12	1.67	0.47
1:C:332:ASN:CG	1:C:333:PRO:HD2	2.39	0.47
1:E:374:HIS:HB3	1:E:377:GLY:O	2.15	0.47
1:H:72:ASP:HA	1:H:73:PRO:HD3	1.83	0.47
1:A:263:PHE:CD2	1:A:314:ARG:HB3	2.50	0.47
1:E:355:GLU:OE2	1:E:393:ARG:NH2	2.49	0.46
1:F:524:ASP:OD2	1:F:524:ASP:N	2.29	0.46
1:E:362:VAL:HG22	1:E:558:LEU:HD21	1.96	0.46
1:G:349:ASP:OD2	1:H:98:LYS:NZ	2.48	0.46
2:I:11:GLY:HA3	2:I:12:GLY:C	2.40	0.46
1:A:95:TYR:CE2	1:A:208:PRO:HD2	2.50	0.46
1:A:374:HIS:HB3	1:A:377:GLY:O	2.16	0.46
1:B:127:LYS:HA	1:B:184:VAL:O	2.16	0.46
1:G:374:HIS:HB3	1:G:377:GLY:O	2.15	0.46
1:H:437:PRO:O	1:H:442:ARG:HD3	2.15	0.46
1:B:488:LEU:HD13	1:B:495:ILE:HD11	1.97	0.46
1:D:420:ARG:NH2	1:D:424:ASP:OD1	2.49	0.46
1:A:139:GLY:HA2	1:A:140:GLY:HA2	1.61	0.46
1:A:362:VAL:HG22	1:A:558:LEU:HD21	1.97	0.46
1:D:142:LEU:HG	1:D:144:PHE:CE1	2.51	0.46
7:G:711:HOH:O	2:J:31:HIS:HE1	1.99	0.46
1:B:243:LYS:HG2	1:B:559:THR:HB	1.98	0.46
1:D:132:ILE:HD12	1:D:180:ASP:HB2	1.98	0.46
1:D:135:LYS:HG3	1:D:145:TYR:CE2	2.50	0.46
1:D:488:LEU:HD13	1:D:495:ILE:HD11	1.96	0.46
1:F:437:PRO:O	1:F:442:ARG:HD3	2.16	0.46
1:A:432:VAL:HG22	1:A:450:GLU:HG3	1.98	0.45
1:C:286:TYR:OH	1:C:291:ASP:HA	2.16	0.45
1:C:307:LEU:HD11	1:D:569:LEU:HD13	1.97	0.45
1:C:385:VAL:HG13	1:C:458:ASN:HA	1.98	0.45
1:H:139:GLY:HA2	7:H:754:HOH:O	2.16	0.45
1:C:121:LEU:O	1:C:187:ASN:HB2	2.15	0.45
1:C:347:TYR:CE1	1:C:484:GLU:HB3	2.51	0.45
1:A:261:LEU:HD13	2:I:20:LEU:HD22	1.97	0.45
1:A:399:GLU:HB3	1:A:452:LEU:HD21	1.97	0.45
1:C:422:ILE:O	1:C:426:ILE:HG13	2.15	0.45
1:E:349:ASP:OD2	1:F:98:LYS:NZ	2.48	0.45
1:G:332:ASN:CG	1:G:333:PRO:HD2	2.42	0.45
1:G:539:TYR:CE1	1:H:232:LEU:HD21	2.52	0.45
1:C:374:HIS:HB3	1:C:377:GLY:O	2.17	0.45
1:D:423:LEU:HD13	1:D:444:LEU:HD13	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:542:PRO:HB3	1:F:181:ILE:HG21	1.99	0.45
1:G:418:GLU:O	1:G:422:ILE:HG13	2.16	0.45
1:A:394:ILE:HB	1:A:399:GLU:HG3	1.97	0.45
1:A:539:TYR:CE1	1:B:232:LEU:HD21	2.50	0.45
1:E:263:PHE:CD2	1:E:314:ARG:HB3	2.52	0.45
1:E:332:ASN:CG	1:E:333:PRO:HD2	2.41	0.45
1:F:127:LYS:HA	1:F:184:VAL:O	2.17	0.45
1:E:394:ILE:HD12	1:E:452:LEU:HD22	1.98	0.45
1:B:361:MET:HE1	1:B:549:MET:HE1	1.99	0.45
7:C:730:HOH:O	1:D:344:TYR:HA	2.17	0.45
1:D:199:ILE:HA	1:D:200:PRO:HD3	1.79	0.45
1:F:423:LEU:HD13	1:F:444:LEU:HD13	1.99	0.45
2:L:13:GLY:HA2	2:L:14:ALA:HA	1.59	0.45
1:G:400:LEU:HD12	1:G:400:LEU:HA	1.80	0.45
1:B:375:PRO:HG3	1:B:383:TYR:CE2	2.51	0.45
1:D:375:PRO:HG3	1:D:383:TYR:CE2	2.52	0.45
1:E:243:LYS:HG2	1:E:559:THR:HB	1.98	0.44
1:H:127:LYS:HA	1:H:184:VAL:O	2.17	0.44
2:K:10:HIS:HB2	2:K:11:GLY:H	1.31	0.44
1:B:135:LYS:HG3	1:B:145:TYR:CE2	2.52	0.44
1:E:457:ILE:HD11	1:E:492:LYS:HE3	1.99	0.44
1:F:375:PRO:HG3	1:F:383:TYR:CE2	2.53	0.44
1:C:157:MET:O	1:C:200:PRO:HD2	2.18	0.44
1:D:392:ARG:NH2	7:D:771:HOH:O	2.49	0.44
1:G:414:PHE:HE2	1:G:444:LEU:HD11	1.82	0.44
1:H:444:LEU:HD12	1:H:444:LEU:HA	1.66	0.44
1:C:191:THR:HB	1:C:193:LYS:NZ	2.32	0.44
1:C:377:GLY:HA2	1:C:378:PRO:HD3	1.84	0.44
1:F:165:SER:HB2	1:F:168:GLU:H	1.82	0.44
1:B:165:SER:HB2	1:B:168:GLU:H	1.83	0.44
1:B:199:ILE:HA	1:B:200:PRO:HD3	1.74	0.44
1:G:362:VAL:HG22	1:G:558:LEU:HD21	1.99	0.44
1:C:108:PHE:CE2	1:C:154:LEU:HD11	2.53	0.44
1:F:372:THR:HG22	1:F:384:ASP:OD2	2.18	0.44
1:A:355:GLU:OE2	1:A:393:ARG:NH2	2.51	0.44
1:E:261:LEU:HD12	1:E:261:LEU:HA	1.85	0.44
1:H:330:THR:O	1:H:331:HIS:CD2	2.71	0.44
1:C:187:ASN:CB	1:C:188:PRO:HD2	2.48	0.44
1:H:164:LYS:HE2	1:H:168:GLU:OE1	2.17	0.44
1:H:372:THR:HG22	1:H:384:ASP:OD2	2.18	0.44
1:H:519:LYS:HE2	1:H:528:PHE:CE2	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:457:ILE:HD11	1:C:492:LYS:HE3	1.99	0.43
1:E:241:ARG:NH2	1:F:311:GLY:O	2.51	0.43
1:E:261:LEU:HD13	2:L:20:LEU:HD22	1.99	0.43
1:E:399:GLU:HB3	1:E:452:LEU:HD21	1.98	0.43
1:A:530:ASP:HB3	1:A:533:PHE:HB3	2.00	0.43
1:E:267:GLU:OE2	1:F:247:ARG:NH2	2.45	0.43
1:E:414:PHE:HE2	1:E:444:LEU:HD11	1.84	0.43
1:E:437:PRO:O	1:E:442:ARG:HD3	2.17	0.43
1:G:394:ILE:HB	1:G:399:GLU:HG3	2.00	0.43
1:B:444:LEU:HD12	1:B:444:LEU:HA	1.66	0.43
1:C:263:PHE:CD2	1:C:314:ARG:HB3	2.53	0.43
1:H:305:LYS:NZ	1:H:501:GLU:OE1	2.37	0.43
1:H:519:LYS:HE2	1:H:528:PHE:CZ	2.53	0.43
1:D:165:SER:HB2	1:D:168:GLU:H	1.83	0.43
1:E:286:TYR:OH	1:E:291:ASP:HA	2.18	0.43
1:H:375:PRO:HG3	1:H:383:TYR:CE2	2.53	0.43
1:D:372:THR:HG22	1:D:384:ASP:OD2	2.19	0.43
1:F:332:ASN:CG	1:F:333:PRO:HD2	2.43	0.43
1:G:399:GLU:HB3	1:G:452:LEU:HD21	2.00	0.43
1:A:568:LEU:O	1:A:571:PRO:HD3	2.19	0.43
1:C:488:LEU:CB	1:C:496:CYS:HB2	2.48	0.43
1:F:72:ASP:HA	1:F:73:PRO:HD3	1.80	0.43
1:F:290:LEU:CD1	1:F:292:MET:HE2	2.49	0.43
1:C:241:ARG:NH1	7:C:715:HOH:O	2.51	0.42
1:C:542:PRO:HB3	1:D:181:ILE:HG21	2.01	0.42
1:E:400:LEU:HD12	1:E:400:LEU:HA	1.82	0.42
1:B:488:LEU:HD23	1:B:488:LEU:HA	1.87	0.42
1:D:125:THR:HG21	1:H:405:GLY:O	2.19	0.42
1:C:254:ILE:HD11	1:C:549:MET:CE	2.50	0.42
1:D:243:LYS:HG2	1:D:559:THR:HB	2.01	0.42
7:F:760:HOH:O	2:L:28:PRO:HD3	2.18	0.42
1:G:98:LYS:NZ	1:H:349:ASP:OD2	2.51	0.42
1:C:394:ILE:HB	1:C:399:GLU:HG3	2.01	0.42
1:D:72:ASP:HA	1:D:73:PRO:HD3	1.80	0.42
1:D:450:GLU:HG2	1:D:451:PHE:CE1	2.54	0.42
1:F:266:ILE:HD13	1:F:307:LEU:CD1	2.50	0.42
1:G:364:HIS:C	1:G:364:HIS:ND1	2.77	0.42
1:H:178:ARG:NH1	1:H:213:LEU:O	2.52	0.42
1:H:213:LEU:HA	1:H:214:PRO:HD3	1.89	0.42
1:B:372:THR:HG22	1:B:384:ASP:OD2	2.19	0.42
1:C:193:LYS:HE3	1:C:193:LYS:N	2.31	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:10:HIS:HB2	2:I:11:GLY:H	1.38	0.42
1:C:329:LEU:HD22	1:C:329:LEU:HA	1.86	0.42
1:F:243:LYS:HG2	1:F:559:THR:HB	2.01	0.42
1:H:135:LYS:HG3	1:H:145:TYR:CE2	2.54	0.42
1:C:270:MET:HE2	1:C:297:ARG:NH2	2.35	0.42
1:C:355:GLU:HA	1:C:391:PHE:CE2	2.55	0.42
1:H:347:TYR:CE1	1:H:484:GLU:HB3	2.55	0.42
2:J:31:HIS:HA	2:J:32:GLY:HA2	1.69	0.42
1:B:420:ARG:NH2	1:B:424:ASP:OD1	2.53	0.42
4:A:602:APC:O3B	4:A:602:APC:O2A	2.38	0.42
1:D:232:LEU:HD23	1:D:232:LEU:HA	1.90	0.42
1:D:122:THR:O	1:H:406:MET:SD	2.77	0.42
1:D:439:THR:OG1	1:D:442:ARG:HG3	2.20	0.42
1:F:373:TYR:CD1	1:F:491:MET:HG3	2.54	0.42
1:C:437:PRO:O	1:C:442:ARG:HD3	2.20	0.41
1:E:220:LEU:HD13	1:E:226:ARG:HG2	2.01	0.41
1:G:488:LEU:CB	1:G:496:CYS:HB2	2.47	0.41
1:A:569:LEU:HD13	1:B:307:LEU:HD11	2.02	0.41
1:B:72:ASP:HA	1:B:73:PRO:HD3	1.81	0.41
1:D:127:LYS:HA	1:D:184:VAL:O	2.20	0.41
1:C:253:TYR:HA	2:K:15:PRO:HB2	2.02	0.41
1:C:270:MET:CE	7:C:739:HOH:O	2.23	0.41
1:E:488:LEU:CB	1:E:496:CYS:HB2	2.50	0.41
1:F:148:ARG:HD3	1:F:153:LYS:HB3	2.02	0.41
2:K:11:GLY:HA3	2:K:12:GLY:C	2.45	0.41
1:G:72:ASP:HA	1:G:73:PRO:HD3	1.88	0.41
1:H:408:LEU:HD23	1:H:408:LEU:HA	1.95	0.41
1:A:108:PHE:CE2	1:A:154:LEU:HD11	2.56	0.41
1:A:484:GLU:HB2	1:A:500:THR:HB	2.03	0.41
1:C:72:ASP:HA	1:C:73:PRO:HD3	1.88	0.41
1:E:300:PRO:HD3	1:E:321:GLN:NE2	2.35	0.41
1:H:243:LYS:HG2	1:H:559:THR:HB	2.02	0.41
1:C:396:MET:HE1	1:C:465:HIS:CD2	2.56	0.41
1:E:266:ILE:HD13	1:E:307:LEU:HD12	2.03	0.41
4:E:602:APC:O2A	4:E:602:APC:O3B	2.39	0.41
1:G:355:GLU:HA	1:G:391:PHE:CE2	2.55	0.41
1:H:423:LEU:HD13	1:H:444:LEU:HD13	2.03	0.41
1:D:308:VAL:HA	1:D:312:ILE:O	2.21	0.41
1:F:100:HIS:HB2	7:F:738:HOH:O	2.21	0.41
1:H:373:TYR:CD1	1:H:491:MET:HG3	2.56	0.41
1:A:300:PRO:HG3	7:A:711:HOH:O	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:404:LEU:HB3	1:A:406:MET:HG2	2.02	0.41
1:A:414:PHE:HE2	1:A:444:LEU:HD11	1.85	0.41
1:C:539:TYR:CE1	1:D:232:LEU:HD21	2.55	0.41
4:C:602:APC:O3B	4:C:602:APC:O2A	2.38	0.41
1:D:266:ILE:HD13	1:D:307:LEU:CD1	2.51	0.41
1:D:309:VAL:HG22	1:D:541:LEU:HA	2.03	0.41
1:E:257:PHE:HA	2:L:16:LEU:CB	2.51	0.41
1:F:199:ILE:HA	1:F:200:PRO:HD3	1.81	0.41
1:G:127:LYS:HA	1:G:184:VAL:O	2.21	0.41
1:G:286:TYR:OH	1:G:291:ASP:HA	2.21	0.41
1:H:199:ILE:HA	1:H:200:PRO:HD3	1.74	0.41
1:H:332:ASN:CG	1:H:333:PRO:HD2	2.46	0.41
2:L:15:PRO:O	2:L:16:LEU:CB	2.69	0.41
1:E:530:ASP:HB3	1:E:533:PHE:HB3	2.02	0.41
1:F:80:ARG:HH12	1:F:180:ASP:CG	2.28	0.41
1:C:414:PHE:HE2	1:C:444:LEU:HD11	1.86	0.40
1:G:305:LYS:NZ	1:G:501:GLU:OE1	2.47	0.40
1:G:569:LEU:HD13	1:H:307:LEU:HD11	2.03	0.40
1:F:439:THR:OG1	1:F:442:ARG:HG3	2.21	0.40
1:F:444:LEU:HD12	1:F:444:LEU:HA	1.69	0.40
1:H:164:LYS:HB2	1:H:164:LYS:HE3	1.90	0.40
1:A:377:GLY:HA2	1:A:378:PRO:HD3	1.83	0.40
1:C:193:LYS:H	1:C:193:LYS:CE	2.30	0.40
1:D:80:ARG:NH1	1:D:180:ASP:OD1	2.55	0.40
1:E:232:LEU:HD23	1:E:232:LEU:HA	1.92	0.40
1:F:139:GLY:HA3	1:F:140:GLY:C	2.46	0.40
1:F:178:ARG:NH1	1:F:213:LEU:O	2.55	0.40
1:G:479:LYS:HD3	1:G:482:LEU:HD12	2.04	0.40
1:G:509:ARG:NH2	1:G:538:GLU:OE2	2.52	0.40
1:H:266:ILE:HD11	1:H:304:HIS:CE1	2.57	0.40
2:L:20:LEU:HD23	2:L:20:LEU:HA	1.97	0.40
1:A:355:GLU:HA	1:A:391:PHE:CE2	2.56	0.40
1:B:266:ILE:HD13	1:B:307:LEU:CD1	2.51	0.40
1:H:172:ILE:HD12	1:H:172:ILE:HA	1.90	0.40
1:A:437:PRO:O	1:A:442:ARG:HD3	2.22	0.40
1:C:432:VAL:HG22	1:C:450:GLU:HG3	2.02	0.40
1:F:512:PHE:HD2	1:F:512:PHE:HA	1.78	0.40
1:H:133:HIS:CD2	1:H:148:ARG:HG3	2.57	0.40
2:K:13:GLY:HA2	2:K:14:ALA:HA	1.67	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	497/513 (97%)	477 (96%)	20 (4%)	0	100	100
1	B	504/513 (98%)	495 (98%)	9 (2%)	0	100	100
1	C	497/513 (97%)	479 (96%)	18 (4%)	0	100	100
1	D	503/513 (98%)	493 (98%)	10 (2%)	0	100	100
1	E	496/513 (97%)	478 (96%)	18 (4%)	0	100	100
1	F	503/513 (98%)	495 (98%)	8 (2%)	0	100	100
1	G	497/513 (97%)	477 (96%)	20 (4%)	0	100	100
1	H	503/513 (98%)	493 (98%)	10 (2%)	0	100	100
2	I	29/54 (54%)	25 (86%)	4 (14%)	0	100	100
2	J	29/54 (54%)	24 (83%)	5 (17%)	0	100	100
2	K	29/54 (54%)	23 (79%)	6 (21%)	0	100	100
2	L	29/54 (54%)	26 (90%)	3 (10%)	0	100	100
All	All	4116/4320 (95%)	3985 (97%)	131 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	409/455 (90%)	374 (91%)	35 (9%)	10	21
1	B	435/455 (96%)	409 (94%)	26 (6%)	17	36

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	425/455 (93%)	387 (91%)	38 (9%)	9	20
1	D	430/455 (94%)	403 (94%)	27 (6%)	16	34
1	E	422/455 (93%)	386 (92%)	36 (8%)	10	22
1	F	430/455 (94%)	402 (94%)	28 (6%)	15	32
1	G	410/455 (90%)	371 (90%)	39 (10%)	8	18
1	H	436/455 (96%)	405 (93%)	31 (7%)	13	29
2	I	21/44 (48%)	18 (86%)	3 (14%)	3	6
2	J	22/44 (50%)	17 (77%)	5 (23%)	1	0
2	K	21/44 (48%)	19 (90%)	2 (10%)	8	18
2	L	23/44 (52%)	19 (83%)	4 (17%)	2	3
All	All	3484/3816 (91%)	3210 (92%)	274 (8%)	11	25

All (274) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	VAL
1	A	74	ASN
1	A	80	ARG
1	A	122	THR
1	A	136	ARG
1	A	138	SER
1	A	143	ILE
1	A	154	LEU
1	A	165	SER
1	A	172	ILE
1	A	212	MET
1	A	220	LEU
1	A	223	LYS
1	A	224	GLU
1	A	240	VAL
1	A	261	LEU
1	A	271	MET
1	A	272	ASN
1	A	328	ASP
1	A	329	LEU
1	A	356	LYS
1	A	365	ILE
1	A	376	ASP

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Mol	Chain	Res	Type
1	A	384	ASP
1	A	404	LEU
1	A	417	GLU
1	A	418	GLU
1	A	444	LEU
1	A	468	ILE
1	A	478	SER
1	A	479	LYS
1	A	488	LEU
1	A	519	LYS
1	A	530	ASP
1	A	549	MET
1	B	75	GLN
1	B	80	ARG
1	B	82	GLN
1	B	133	HIS
1	B	147	LEU
1	B	152	VAL
1	B	175	LYS
1	B	207	SER
1	B	220	LEU
1	B	266	ILE
1	B	271	MET
1	B	272	ASN
1	B	274	ILE
1	B	279	VAL
1	B	330	THR
1	B	347	TYR
1	B	355	GLU
1	B	364	HIS
1	B	371	VAL
1	B	404	LEU
1	B	415	GLU
1	B	420	ARG
1	B	488	LEU
1	B	524	ASP
1	B	549	MET
1	B	552	ASP
1	C	74	ASN
1	C	80	ARG
1	C	122	THR
1	C	136	ARG

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Mol	Chain	Res	Type
1	C	138	SER
1	C	141	LYS
1	C	154	LEU
1	C	165	SER
1	C	168	GLU
1	C	172	ILE
1	C	187	ASN
1	C	193	LYS
1	C	212	MET
1	C	220	LEU
1	C	223	LYS
1	C	224	GLU
1	C	240	VAL
1	C	260	GLU
1	C	261	LEU
1	C	271	MET
1	C	272	ASN
1	C	279	VAL
1	C	328	ASP
1	C	329	LEU
1	C	347	TYR
1	C	352	GLU
1	C	365	ILE
1	C	376	ASP
1	C	384	ASP
1	C	404	LEU
1	C	418	GLU
1	C	444	LEU
1	C	468	ILE
1	C	478	SER
1	C	479	LYS
1	C	488	LEU
1	C	530	ASP
1	C	549	MET
1	D	75	GLN
1	D	80	ARG
1	D	82	GLN
1	D	133	HIS
1	D	138	SER
1	D	147	LEU
1	D	152	VAL
1	D	175	LYS

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Mol	Chain	Res	Type
1	D	220	LEU
1	D	266	ILE
1	D	271	MET
1	D	272	ASN
1	D	274	ILE
1	D	279	VAL
1	D	320	ARG
1	D	330	THR
1	D	347	TYR
1	D	355	GLU
1	D	363	LYS
1	D	371	VAL
1	D	393	ARG
1	D	404	LEU
1	D	415	GLU
1	D	420	ARG
1	D	488	LEU
1	D	549	MET
1	D	552	ASP
1	E	74	ASN
1	E	75	GLN
1	E	80	ARG
1	E	82	GLN
1	E	111	LYS
1	E	116	GLN
1	E	122	THR
1	E	136	ARG
1	E	138	SER
1	E	154	LEU
1	E	165	SER
1	E	172	ILE
1	E	187	ASN
1	E	212	MET
1	E	220	LEU
1	E	224	GLU
1	E	240	VAL
1	E	260	GLU
1	E	261	LEU
1	E	271	MET
1	E	272	ASN
1	E	327	ILE
1	E	328	ASP

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Mol	Chain	Res	Type
1	E	347	TYR
1	E	352	GLU
1	E	365	ILE
1	E	376	ASP
1	E	379	GLU
1	E	393	ARG
1	E	404	LEU
1	E	444	LEU
1	E	468	ILE
1	E	479	LYS
1	E	488	LEU
1	E	530	ASP
1	E	549	MET
1	F	71	VAL
1	F	75	GLN
1	F	80	ARG
1	F	82	GLN
1	F	133	HIS
1	F	147	LEU
1	F	152	VAL
1	F	175	LYS
1	F	207	SER
1	F	220	LEU
1	F	266	ILE
1	F	271	MET
1	F	272	ASN
1	F	274	ILE
1	F	279	VAL
1	F	330	THR
1	F	347	TYR
1	F	355	GLU
1	F	363	LYS
1	F	370	LYS
1	F	371	VAL
1	F	404	LEU
1	F	415	GLU
1	F	488	LEU
1	F	524	ASP
1	F	549	MET
1	F	552	ASP
1	F	574	LYS
1	G	74	ASN

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Mol	Chain	Res	Type
1	G	80	ARG
1	G	122	THR
1	G	136	ARG
1	G	138	SER
1	G	154	LEU
1	G	165	SER
1	G	172	ILE
1	G	187	ASN
1	G	212	MET
1	G	220	LEU
1	G	223	LYS
1	G	224	GLU
1	G	240	VAL
1	G	261	LEU
1	G	271	MET
1	G	272	ASN
1	G	279	VAL
1	G	284	ILE
1	G	321	GLN
1	G	327	ILE
1	G	347	TYR
1	G	352	GLU
1	G	365	ILE
1	G	376	ASP
1	G	384	ASP
1	G	392	ARG
1	G	393	ARG
1	G	402	LYS
1	G	404	LEU
1	G	407	LYS
1	G	417	GLU
1	G	427	CYS
1	G	444	LEU
1	G	468	ILE
1	G	479	LYS
1	G	488	LEU
1	G	530	ASP
1	G	549	MET
1	H	71	VAL
1	H	75	GLN
1	H	80	ARG
1	H	82	GLN

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Mol	Chain	Res	Type
1	H	133	HIS
1	H	142	LEU
1	H	147	LEU
1	H	152	VAL
1	H	175	LYS
1	H	207	SER
1	H	220	LEU
1	H	221	LYS
1	H	266	ILE
1	H	271	MET
1	H	272	ASN
1	H	274	ILE
1	H	279	VAL
1	H	327	ILE
1	H	328	ASP
1	H	330	THR
1	H	347	TYR
1	H	355	GLU
1	H	363	LYS
1	H	370	LYS
1	H	371	VAL
1	H	404	LEU
1	H	470	SER
1	H	488	LEU
1	H	524	ASP
1	H	549	MET
1	H	552	ASP
2	I	5	GLN
2	I	20	LEU
2	I	30	VAL
2	J	5	GLN
2	J	10	HIS
2	J	18	VAL
2	J	20	LEU
2	J	30	VAL
2	K	5	GLN
2	K	30	VAL
2	L	5	GLN
2	L	10	HIS
2	L	18	VAL
2	L	30	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	HIS
1	A	229	GLN
1	A	272	ASN
1	A	288	ASN
1	A	321	GLN
1	A	364	HIS
1	B	185	GLN
1	B	229	GLN
1	B	348	HIS
1	C	85	HIS
1	C	116	GLN
1	C	288	ASN
1	C	348	HIS
1	C	364	HIS
1	D	75	GLN
1	D	229	GLN
1	D	321	GLN
1	D	348	HIS
1	D	563	ASN
1	E	85	HIS
1	E	86	GLN
1	E	90	ASN
1	E	187	ASN
1	E	229	GLN
1	E	272	ASN
1	E	287	HIS
1	E	288	ASN
1	E	348	HIS
1	F	75	GLN
1	F	229	GLN
1	F	321	GLN
1	F	348	HIS
1	G	85	HIS
1	G	187	ASN
1	G	229	GLN
1	G	272	ASN
1	G	287	HIS
1	G	288	ASN
1	G	348	HIS
1	H	229	GLN
1	H	288	ASN
1	H	321	GLN

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Mol	Chain	Res	Type
1	H	331	HIS
1	H	348	HIS
2	I	31	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 29 ligands modelled in this entry, 12 are monoatomic - leaving 17 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	APC	G	602	-	26,29,33	1.39	5 (19%)	40,45,52	1.87	9 (22%)
3	LYS	G	601	-	8,9,9	0.84	1 (12%)	7,10,10	1.00	1 (14%)
4	APC	F	602	5	29,33,33	1.46	5 (17%)	44,52,52	1.84	11 (25%)
4	APC	D	602	5	29,33,33	1.46	6 (20%)	44,52,52	1.84	12 (27%)
6	ALA	H	601	-	3,4,5	0.96	0	2,4,6	0.55	0
3	LYS	B	601	-	8,9,9	0.79	1 (12%)	7,10,10	0.94	1 (14%)
3	LYS	A	601	-	8,9,9	0.77	0	7,10,10	1.11	1 (14%)
4	APC	E	602	-	26,29,33	1.34	5 (19%)	40,45,52	1.94	9 (22%)
4	APC	A	602	-	26,29,33	1.33	5 (19%)	40,45,52	1.89	9 (22%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	APC	C	602	-	26,29,33	1.32	5 (19%)	40,45,52	1.93	9 (22%)
3	LYS	E	601	-	8,9,9	0.85	1 (12%)	7,10,10	1.31	1 (14%)
4	APC	H	603	5	29,33,33	1.47	5 (17%)	44,52,52	1.82	12 (27%)
3	LYS	F	601	-	8,9,9	0.82	0	7,10,10	1.11	1 (14%)
3	LYS	H	602	-	8,9,9	0.85	1 (12%)	7,10,10	1.10	1 (14%)
4	APC	B	602	5	29,33,33	1.45	6 (20%)	44,52,52	1.81	10 (22%)
3	LYS	C	601	-	8,9,9	0.79	0	7,10,10	1.29	1 (14%)
3	LYS	D	601	-	8,9,9	0.89	1 (12%)	7,10,10	1.16	1 (14%)

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	APC	G	602	-	-	9/16/32/38	0/3/3/3
3	LYS	G	601	-	-	3/9/9/9	-
4	APC	F	602	5	-	3/19/38/38	0/3/3/3
4	APC	D	602	5	-	3/19/38/38	0/3/3/3
6	ALA	H	601	-	-	0/1/2/4	-
3	LYS	B	601	-	-	2/9/9/9	-
3	LYS	A	601	-	-	3/9/9/9	-
4	APC	E	602	-	-	9/16/32/38	0/3/3/3
4	APC	A	602	-	-	9/16/32/38	0/3/3/3
4	APC	C	602	-	-	9/16/32/38	0/3/3/3
3	LYS	E	601	-	-	3/9/9/9	-
4	APC	H	603	5	-	3/19/38/38	0/3/3/3
3	LYS	F	601	-	-	1/9/9/9	-
3	LYS	H	602	-	-	0/9/9/9	-
4	APC	B	602	5	-	3/19/38/38	0/3/3/3
3	LYS	C	601	-	-	3/9/9/9	-
3	LYS	D	601	-	-	1/9/9/9	-

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	602	APC	PG-O1G	3.63	1.61	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	602	APC	PG-O1G	3.57	1.61	1.50
4	F	602	APC	PG-O1G	3.56	1.61	1.50
4	H	603	APC	PG-O1G	3.50	1.61	1.50
4	B	602	APC	PB-O3B	3.42	1.62	1.58
4	F	602	APC	PB-O3B	3.36	1.62	1.58
4	G	602	APC	PB-O3B	3.33	1.62	1.55
4	A	602	APC	PB-O3B	3.30	1.62	1.55
4	D	602	APC	PB-O3B	3.26	1.62	1.58
4	H	603	APC	PB-O3B	3.24	1.62	1.58
4	G	602	APC	PB-O2B	3.22	1.62	1.55
4	E	602	APC	PB-O3B	3.11	1.61	1.55
4	C	602	APC	PB-O3B	2.98	1.61	1.55
4	A	602	APC	PB-O2B	2.92	1.61	1.55
4	C	602	APC	PB-O2B	2.92	1.61	1.55
4	E	602	APC	PB-O2B	2.85	1.61	1.55
4	H	603	APC	C8-N7	2.33	1.36	1.31
4	C	602	APC	PA-O2A	2.30	1.62	1.56
4	G	602	APC	PA-O2A	2.28	1.62	1.56
4	A	602	APC	PA-O2A	2.28	1.62	1.56
4	G	602	APC	C5-N7	-2.27	1.34	1.39
4	E	602	APC	PA-O2A	2.26	1.61	1.56
4	H	603	APC	C5-N7	-2.22	1.35	1.39
4	F	602	APC	C8-N7	2.22	1.35	1.31
4	D	602	APC	C8-N7	2.18	1.35	1.31
4	D	602	APC	PB-O2B	2.18	1.61	1.56
4	G	602	APC	C8-N7	2.17	1.35	1.31
4	E	602	APC	C8-N7	2.17	1.35	1.31
4	D	602	APC	C5-N7	-2.16	1.35	1.39
4	E	602	APC	C5-N7	-2.14	1.35	1.39
4	F	602	APC	C5-N7	-2.12	1.35	1.39
3	E	601	LYS	OXT-C	-2.11	1.23	1.30
4	A	602	APC	C5-N7	-2.11	1.35	1.39
4	F	602	APC	PB-O2B	2.09	1.61	1.56
4	B	602	APC	PB-O2B	2.09	1.61	1.56
3	D	601	LYS	OXT-C	-2.08	1.24	1.30
3	G	601	LYS	OXT-C	-2.08	1.24	1.30
4	A	602	APC	C8-N7	2.07	1.35	1.31
4	C	602	APC	C5-N7	-2.07	1.35	1.39
3	H	602	LYS	OXT-C	-2.05	1.24	1.30
4	B	602	APC	C8-N7	2.05	1.35	1.31
4	B	602	APC	C5-N7	-2.04	1.35	1.39
4	H	603	APC	PA-O2A	2.04	1.61	1.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	601	LYS	OXT-C	-2.03	1.24	1.30
4	C	602	APC	C8-N7	2.03	1.35	1.31
4	B	602	APC	PA-O2A	2.01	1.61	1.56
4	D	602	APC	PA-O2A	2.00	1.61	1.56

All (89) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	602	APC	C5-C4-N3	-5.45	119.21	126.72
4	C	602	APC	C5-C4-N3	-5.34	119.36	126.72
4	D	602	APC	C5-C4-N3	-5.31	119.40	126.72
4	A	602	APC	C5-C4-N3	-5.31	119.41	126.72
4	G	602	APC	C5-C4-N3	-5.25	119.48	126.72
4	B	602	APC	C5-C4-N3	-5.10	119.69	126.72
4	F	602	APC	C5-C4-N3	-5.10	119.70	126.72
4	C	602	APC	N3-C2-N1	-5.05	120.94	128.58
4	E	602	APC	N3-C2-N1	-5.04	120.96	128.58
4	A	602	APC	N3-C2-N1	-4.94	121.11	128.58
4	G	602	APC	N3-C2-N1	-4.93	121.12	128.58
4	H	603	APC	C5-C4-N3	-4.91	119.96	126.72
4	H	603	APC	N3-C2-N1	-4.89	121.17	128.58
4	F	602	APC	N3-C2-N1	-4.85	121.24	128.58
4	B	602	APC	N3-C2-N1	-4.55	121.69	128.58
4	D	602	APC	N3-C2-N1	-4.45	121.85	128.58
4	E	602	APC	N3-C4-N9	3.68	133.43	127.17
4	C	602	APC	N3-C4-N9	3.61	133.31	127.17
4	A	602	APC	N3-C4-N9	3.58	133.26	127.17
4	G	602	APC	N3-C4-N9	3.55	133.21	127.17
4	E	602	APC	C2-N3-C4	3.50	120.39	111.83
4	C	602	APC	C2-N3-C4	3.43	120.20	111.83
4	F	602	APC	N3-C4-N9	3.42	132.98	127.17
4	D	602	APC	N3-C4-N9	3.41	132.97	127.17
4	A	602	APC	C2-N3-C4	3.41	120.15	111.83
4	G	602	APC	C2-N3-C4	3.40	120.13	111.83
4	H	603	APC	N9-C8-N7	-3.39	109.12	113.94
4	B	602	APC	N9-C8-N7	-3.39	109.12	113.94
4	D	602	APC	N9-C8-N7	-3.38	109.14	113.94
3	E	601	LYS	OXT-C-O	-3.37	116.44	124.08
4	C	602	APC	N9-C8-N7	-3.36	109.18	113.94
4	B	602	APC	N3-C4-N9	3.34	132.85	127.17
4	F	602	APC	N9-C8-N7	-3.33	109.22	113.94
3	C	601	LYS	OXT-C-O	-3.29	116.61	124.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	602	APC	N9-C8-N7	-3.29	109.28	113.94
4	E	602	APC	N9-C8-N7	-3.27	109.30	113.94
4	H	603	APC	N3-C4-N9	3.26	132.71	127.17
4	F	602	APC	C2-N3-C4	3.26	119.79	111.83
4	D	602	APC	C2-N3-C4	3.21	119.68	111.83
4	B	602	APC	C2-N3-C4	3.16	119.55	111.83
4	H	603	APC	C2-N3-C4	3.10	119.41	111.83
4	D	602	APC	C4-C5-N7	-3.06	107.08	110.58
4	G	602	APC	N9-C8-N7	-3.04	109.63	113.94
4	B	602	APC	C4-C5-N7	-2.98	107.17	110.58
3	D	601	LYS	OXT-C-O	-2.94	117.40	124.08
4	H	603	APC	PB-O3B-PG	-2.93	121.97	132.45
4	D	602	APC	PB-O3B-PG	-2.89	122.11	132.45
4	D	602	APC	C5-N7-C8	2.88	107.98	103.45
4	F	602	APC	PB-O3B-PG	-2.86	122.20	132.45
4	B	602	APC	C5-N7-C8	2.86	107.94	103.45
4	C	602	APC	C4-C5-N7	-2.84	107.33	110.58
4	B	602	APC	PB-O3B-PG	-2.84	122.27	132.45
3	A	601	LYS	OXT-C-O	-2.82	117.69	124.08
3	H	602	LYS	OXT-C-O	-2.79	117.74	124.08
4	C	602	APC	C5-N7-C8	2.78	107.82	103.45
3	F	601	LYS	OXT-C-O	-2.76	117.82	124.08
4	F	602	APC	C4-C5-N7	-2.74	107.45	110.58
4	H	603	APC	C5-N7-C8	2.74	107.75	103.45
4	E	602	APC	C4-C5-N7	-2.73	107.46	110.58
4	E	602	APC	C5-N7-C8	2.73	107.74	103.45
4	F	602	APC	C5-N7-C8	2.73	107.74	103.45
4	H	603	APC	C4-C5-N7	-2.72	107.47	110.58
4	A	602	APC	C5-N7-C8	2.69	107.67	103.45
4	A	602	APC	C4-C5-N7	-2.67	107.53	110.58
4	G	602	APC	C5-N7-C8	2.51	107.40	103.45
3	G	601	LYS	OXT-C-O	-2.49	118.44	124.08
4	C	602	APC	C4-N9-C8	2.48	108.34	105.74
4	H	603	APC	C4-N9-C8	2.47	108.33	105.74
4	G	602	APC	C4-C5-N7	-2.46	107.77	110.58
4	B	602	APC	C4-N9-C8	2.45	108.31	105.74
4	E	602	APC	C3'-C2'-C1'	2.43	106.06	101.46
4	G	602	APC	C3'-C2'-C1'	2.43	106.06	101.46
4	A	602	APC	C3'-C2'-C1'	2.43	106.05	101.46
4	F	602	APC	C4-N9-C8	2.42	108.28	105.74
4	A	602	APC	C4-N9-C8	2.34	108.19	105.74
4	C	602	APC	C3'-C2'-C1'	2.32	105.85	101.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	602	APC	C4-N9-C8	2.31	108.17	105.74
3	B	601	LYS	OXT-C-O	-2.30	118.86	124.08
4	E	602	APC	C4-N9-C8	2.29	108.14	105.74
4	H	603	APC	C3'-C2'-C1'	2.18	105.59	101.46
4	B	602	APC	C2'-C3'-C4'	2.17	106.79	102.61
4	F	602	APC	C3'-C2'-C1'	2.14	105.52	101.46
4	D	602	APC	C3'-C2'-C1'	2.13	105.50	101.46
4	H	603	APC	C2-N1-C6	2.13	122.23	118.73
4	F	602	APC	C2'-C3'-C4'	2.12	106.70	102.61
4	G	602	APC	C4-N9-C8	2.06	107.90	105.74
4	H	603	APC	C2'-C3'-C4'	2.05	106.56	102.61
4	D	602	APC	C2'-C3'-C4'	2.04	106.56	102.61
4	D	602	APC	C6-C5-C4	2.04	119.96	117.18

There are no chirality outliers.

All (64) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	602	APC	PA-C3A-PB-O1B
4	A	602	APC	PA-C3A-PB-O2B
4	A	602	APC	PA-C3A-PB-O3B
4	A	602	APC	C5'-O5'-PA-O1A
4	A	602	APC	C5'-O5'-PA-C3A
4	A	602	APC	O4'-C4'-C5'-O5'
4	B	602	APC	C5'-O5'-PA-O1A
4	C	602	APC	PA-C3A-PB-O1B
4	C	602	APC	PA-C3A-PB-O2B
4	C	602	APC	PA-C3A-PB-O3B
4	C	602	APC	C5'-O5'-PA-O1A
4	C	602	APC	C5'-O5'-PA-C3A
4	C	602	APC	O4'-C4'-C5'-O5'
4	D	602	APC	C5'-O5'-PA-O1A
4	E	602	APC	PA-C3A-PB-O1B
4	E	602	APC	PA-C3A-PB-O2B
4	E	602	APC	PA-C3A-PB-O3B
4	E	602	APC	C5'-O5'-PA-O1A
4	E	602	APC	C5'-O5'-PA-C3A
4	E	602	APC	O4'-C4'-C5'-O5'
4	F	602	APC	C5'-O5'-PA-O1A
4	G	602	APC	PA-C3A-PB-O1B
4	G	602	APC	PA-C3A-PB-O2B
4	G	602	APC	PA-C3A-PB-O3B

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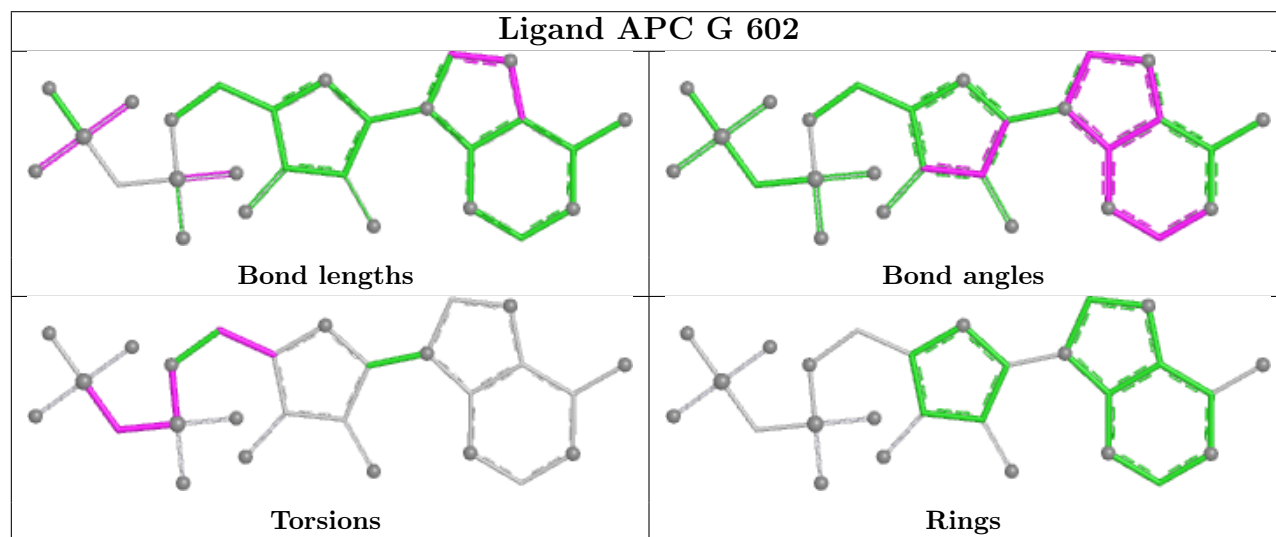
Mol	Chain	Res	Type	Atoms
4	G	602	APC	C5'-O5'-PA-O1A
4	G	602	APC	C5'-O5'-PA-C3A
4	G	602	APC	O4'-C4'-C5'-O5'
4	H	603	APC	C5'-O5'-PA-O1A
4	A	602	APC	C3'-C4'-C5'-O5'
4	C	602	APC	C3'-C4'-C5'-O5'
4	E	602	APC	C3'-C4'-C5'-O5'
4	G	602	APC	C3'-C4'-C5'-O5'
4	H	603	APC	O4'-C4'-C5'-O5'
4	B	602	APC	O4'-C4'-C5'-O5'
4	F	602	APC	O4'-C4'-C5'-O5'
3	C	601	LYS	CA-CB-CG-CD
3	A	601	LYS	CA-CB-CG-CD
4	D	602	APC	O4'-C4'-C5'-O5'
4	A	602	APC	PB-C3A-PA-O2A
4	C	602	APC	PB-C3A-PA-O2A
4	E	602	APC	PB-C3A-PA-O2A
4	G	602	APC	PB-C3A-PA-O2A
3	E	601	LYS	CA-CB-CG-CD
3	G	601	LYS	CA-CB-CG-CD
3	B	601	LYS	O-C-CA-N
3	D	601	LYS	O-C-CA-N
3	F	601	LYS	O-C-CA-N
3	A	601	LYS	O-C-CA-CB
3	E	601	LYS	O-C-CA-CB
3	E	601	LYS	OXT-C-CA-CB
3	G	601	LYS	O-C-CA-CB
3	G	601	LYS	OXT-C-CA-CB
4	A	602	APC	PB-C3A-PA-O1A
4	B	602	APC	C5'-O5'-PA-O2A
4	C	602	APC	PB-C3A-PA-O1A
4	D	602	APC	C5'-O5'-PA-O2A
4	E	602	APC	PB-C3A-PA-O1A
4	F	602	APC	C5'-O5'-PA-O2A
4	G	602	APC	PB-C3A-PA-O1A
4	H	603	APC	C5'-O5'-PA-O2A
3	A	601	LYS	OXT-C-CA-CB
3	B	601	LYS	CA-CB-CG-CD
3	C	601	LYS	O-C-CA-CB
3	C	601	LYS	OXT-C-CA-CB

There are no ring outliers.

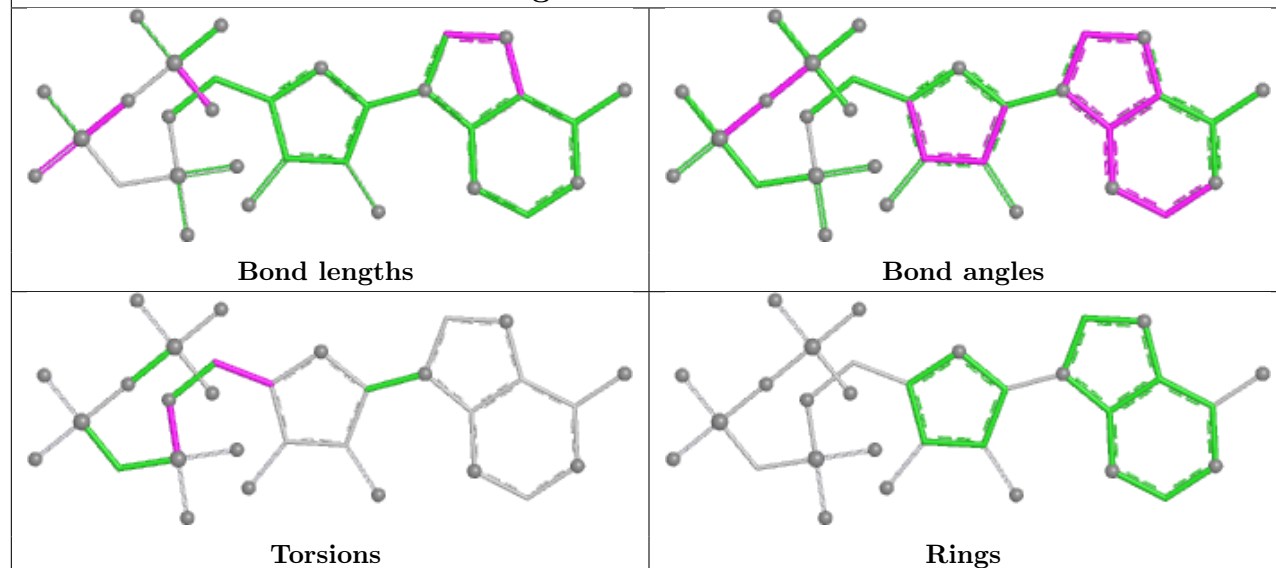
4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	H	601	ALA	1	0
4	E	602	APC	2	0
4	A	602	APC	1	0
4	C	602	APC	1	0

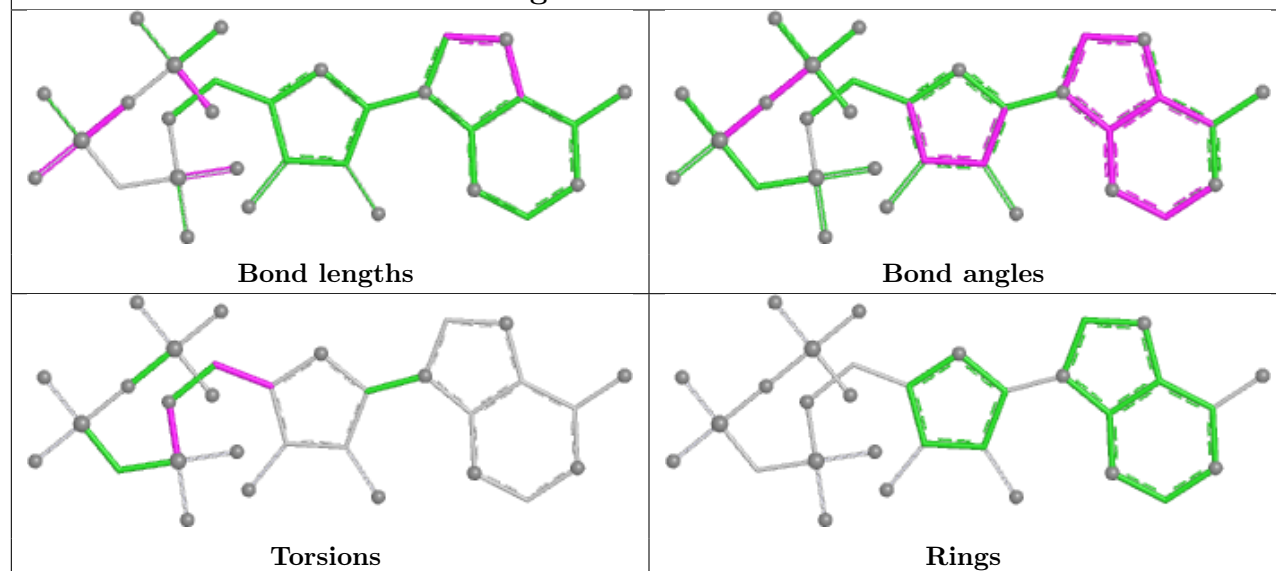
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



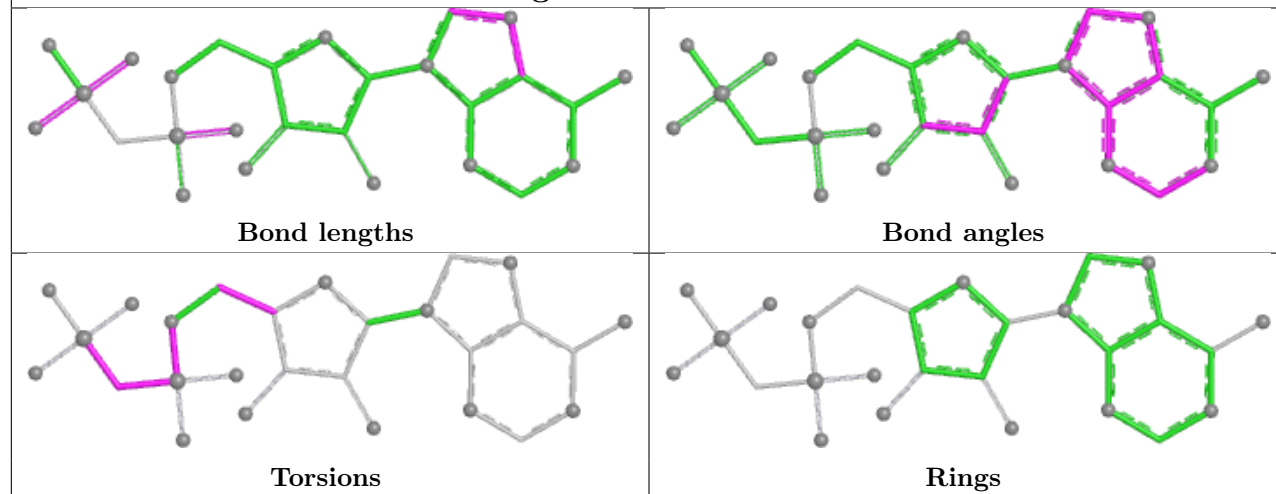
Ligand APC F 602

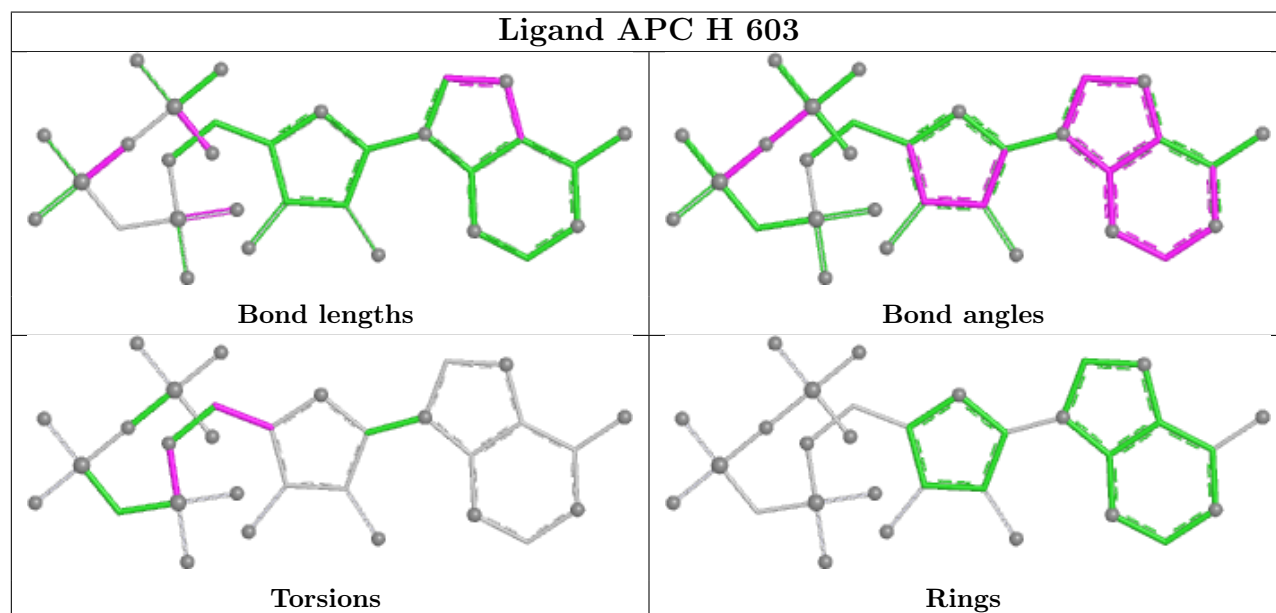
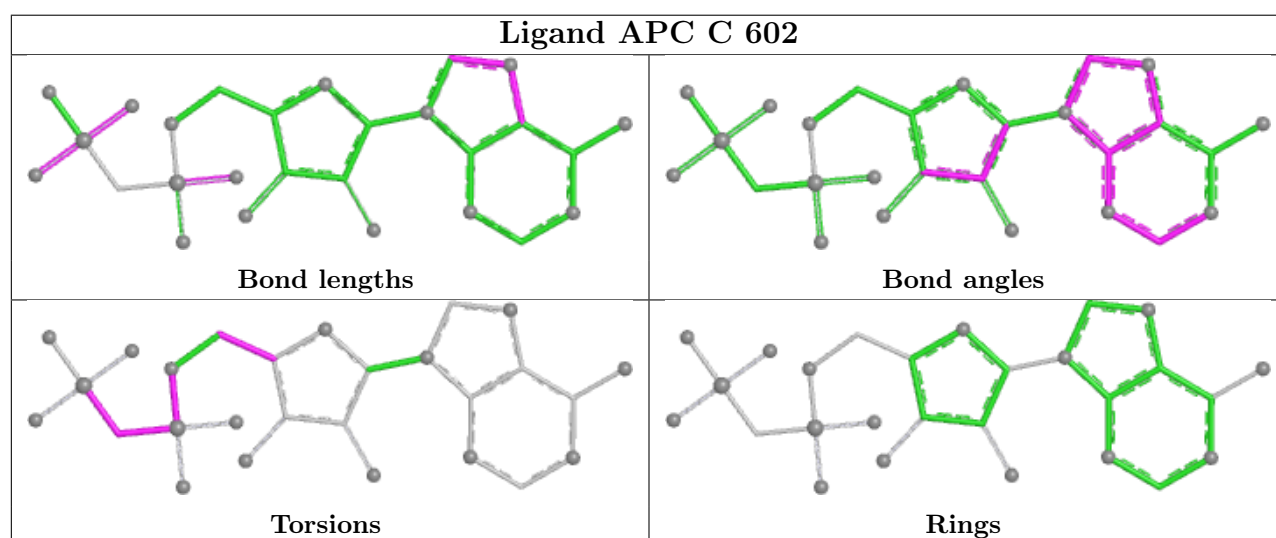
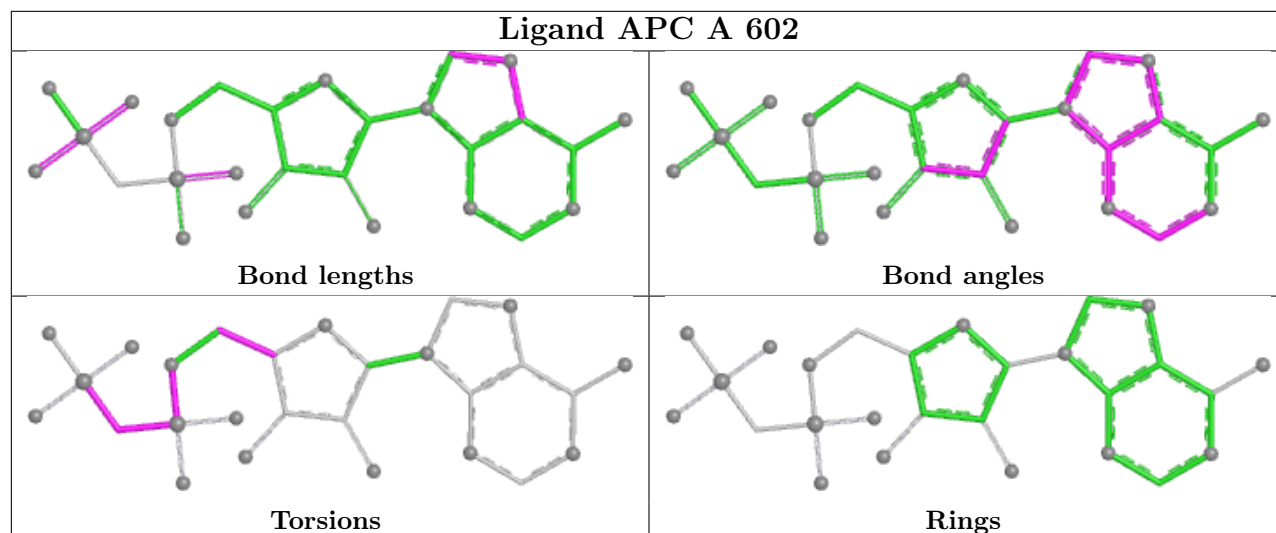


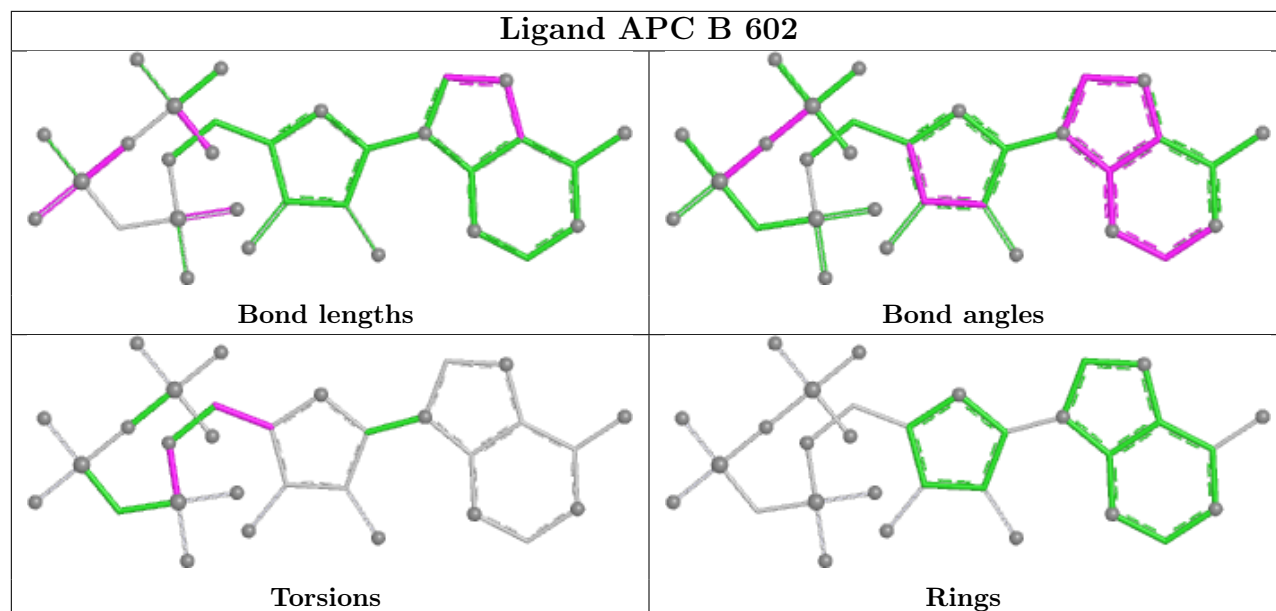
Ligand APC D 602



Ligand APC E 602







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	501/513 (97%)	-1.71	0 100 100	17, 49, 84, 100	0
1	B	506/513 (98%)	-1.85	0 100 100	17, 33, 66, 82	0
1	C	501/513 (97%)	-1.77	0 100 100	16, 40, 74, 97	0
1	D	505/513 (98%)	-1.82	0 100 100	16, 34, 70, 112	0
1	E	500/513 (97%)	-1.79	0 100 100	16, 39, 75, 114	0
1	F	505/513 (98%)	-1.83	0 100 100	16, 32, 72, 117	0
1	G	501/513 (97%)	-1.70	0 100 100	19, 49, 83, 108	0
1	H	505/513 (98%)	-1.85	0 100 100	18, 33, 65, 91	0
2	I	31/54 (57%)	-1.40	0 100 100	23, 50, 114, 118	0
2	J	31/54 (57%)	-1.42	0 100 100	21, 53, 115, 125	0
2	K	31/54 (57%)	-1.34	0 100 100	21, 53, 113, 118	0
2	L	31/54 (57%)	-1.38	0 100 100	20, 49, 114, 116	0
All	All	4148/4320 (96%)	-1.78	0 100 100	16, 37, 79, 125	0

There are no RSRZ outliers to report.

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 ⓘ

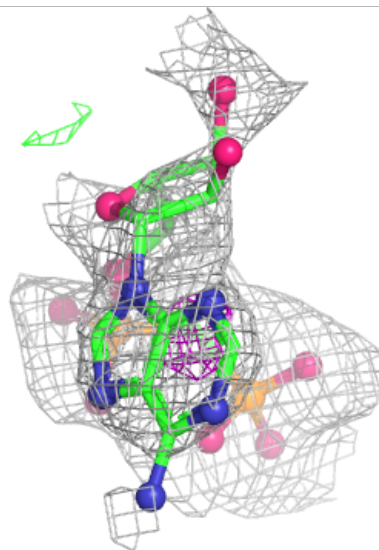
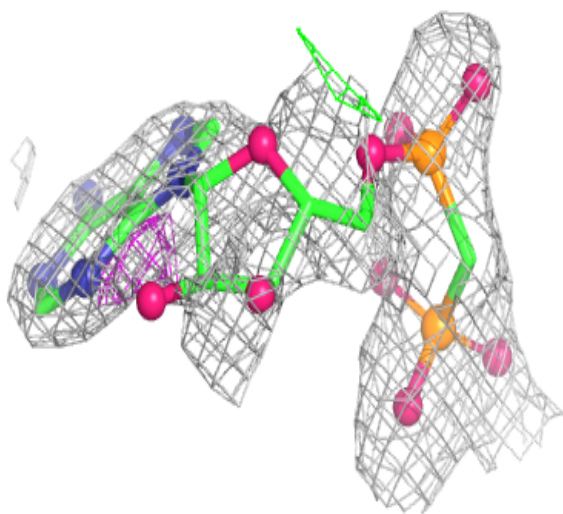
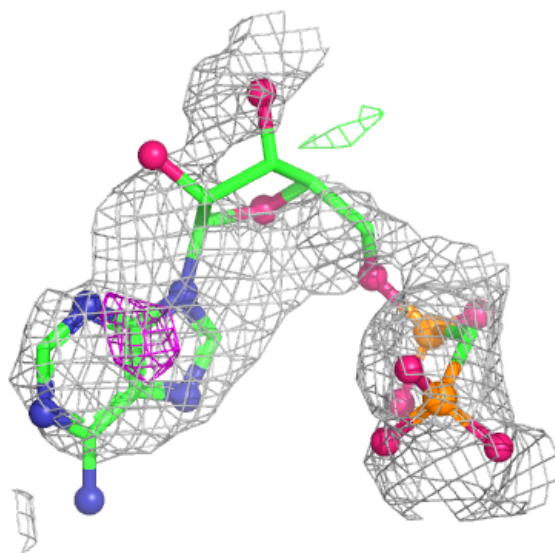
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	ALA	H	601	5/6	0.95	0.09	69,76,79,80	0
5	MG	G	603	1/1	0.98	0.03	55,55,55,55	0
4	APC	C	602	27/31	0.98	0.06	72,101,131,155	0
4	APC	D	602	31/31	0.99	0.05	60,108,149,152	0
4	APC	E	602	27/31	0.99	0.05	74,99,127,160	0
4	APC	F	602	31/31	0.99	0.05	61,107,145,155	0
4	APC	G	602	27/31	0.99	0.06	83,111,142,148	0
4	APC	H	603	31/31	0.99	0.05	56,105,139,158	0
5	MG	A	603	1/1	0.99	0.02	53,53,53,53	0
5	MG	B	603	1/1	0.99	0.02	63,63,63,63	0
5	MG	C	603	1/1	0.99	0.03	59,59,59,59	0
5	MG	E	603	1/1	0.99	0.03	53,53,53,53	0
4	APC	B	602	31/31	0.99	0.04	57,101,132,156	0
5	MG	H	604	1/1	0.99	0.03	66,66,66,66	0
4	APC	A	602	27/31	0.99	0.06	73,101,138,154	0
3	LYS	H	602	10/10	1.00	0.02	18,22,30,32	0
3	LYS	A	601	10/10	1.00	0.03	16,24,46,61	0
3	LYS	B	601	10/10	1.00	0.02	15,24,32,35	0
3	LYS	C	601	10/10	1.00	0.02	15,19,42,44	0
5	MG	D	603	1/1	1.00	0.02	68,68,68,68	0
3	LYS	D	601	10/10	1.00	0.02	15,31,36,38	0
5	MG	F	603	1/1	1.00	0.01	59,59,59,59	0
5	MG	F	604	1/1	1.00	0.03	12,12,12,12	0
3	LYS	E	601	10/10	1.00	0.02	15,22,45,49	0
3	LYS	F	601	10/10	1.00	0.02	17,30,36,39	0
5	MG	H	605	1/1	1.00	0.01	13,13,13,13	0
5	MG	J	101	1/1	1.00	0.02	21,21,21,21	0
5	MG	L	101	1/1	1.00	0.02	20,20,20,20	0
3	LYS	G	601	10/10	1.00	0.03	17,33,43,52	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.

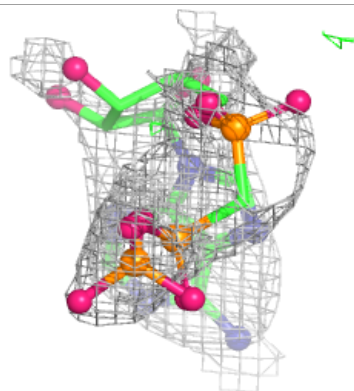
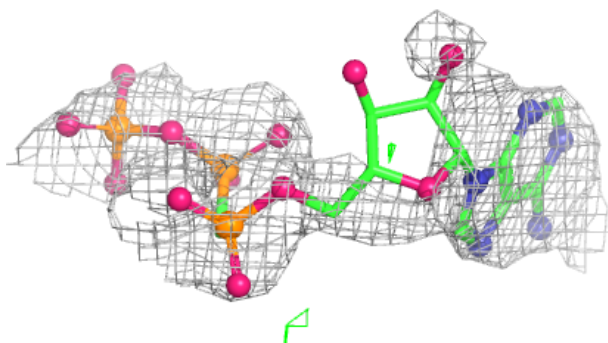
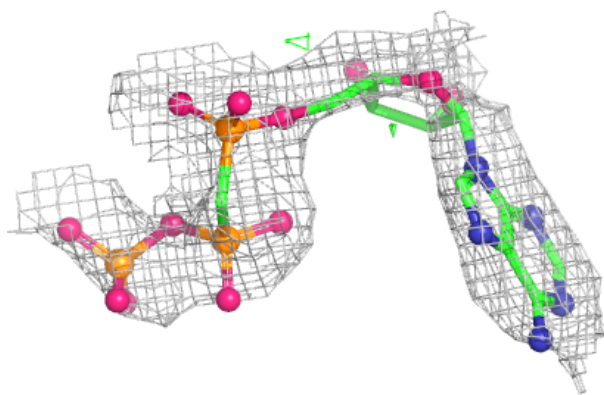
Electron density around APC C 602:

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



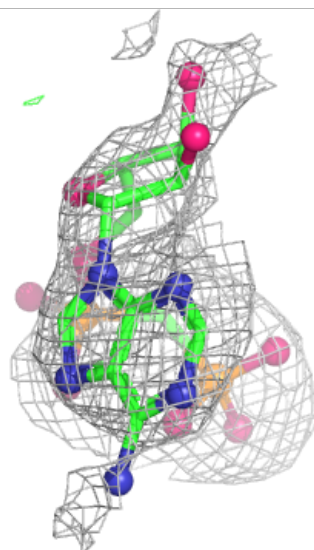
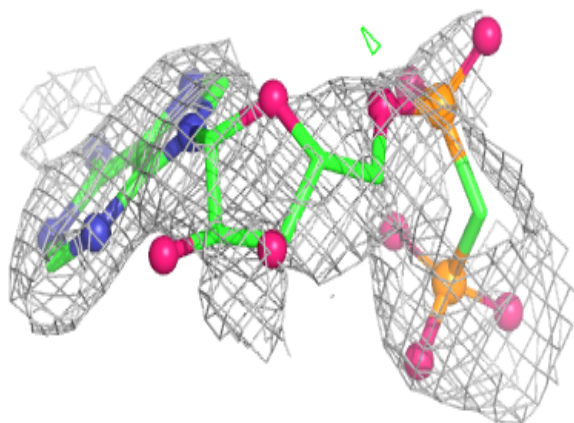
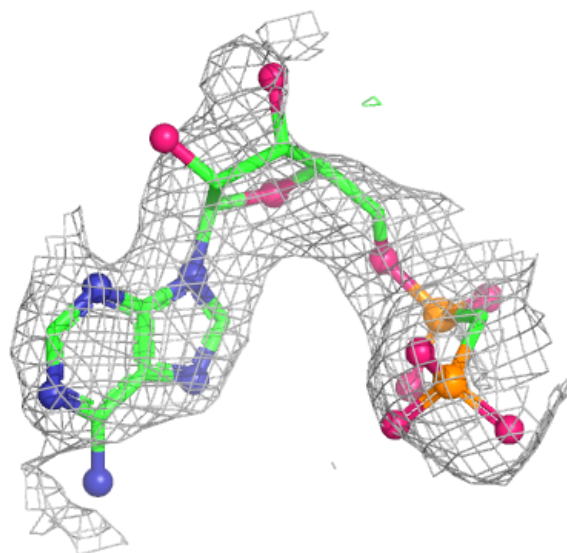
Electron density around APC D 602:

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



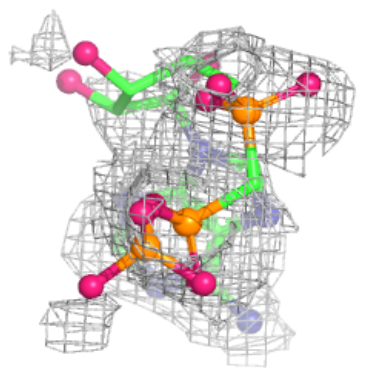
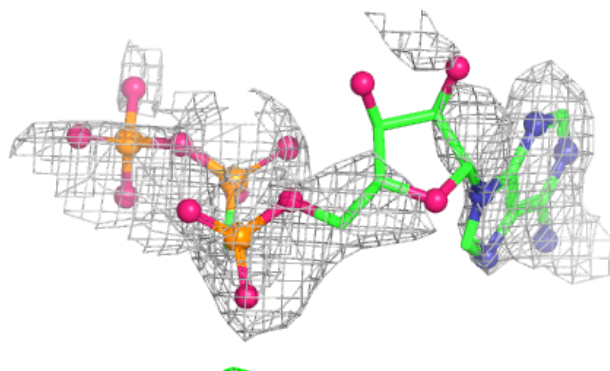
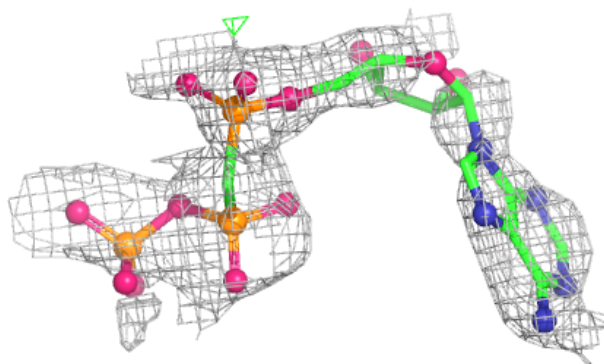
Electron density around APC E 602:

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



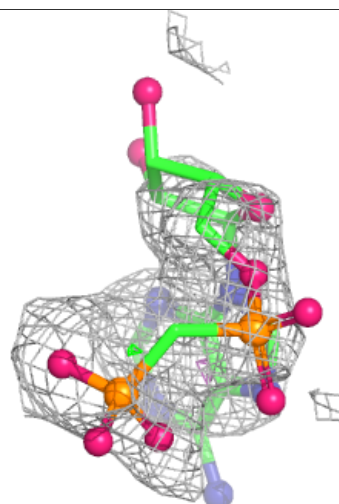
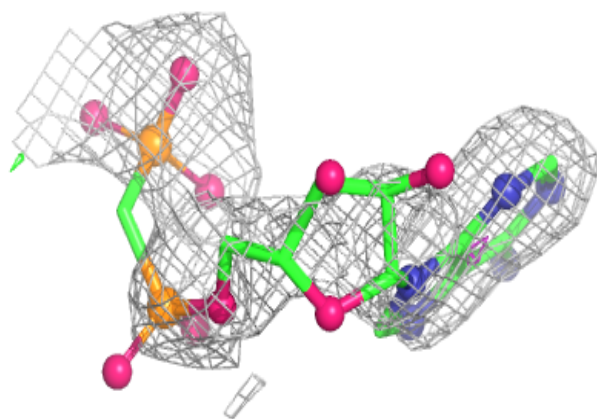
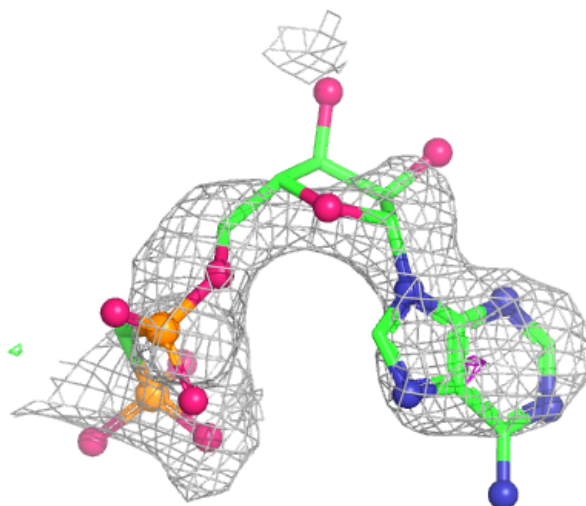
Electron density around APC F 602:

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



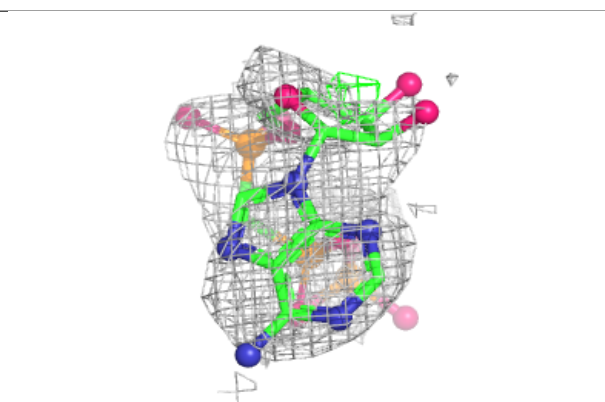
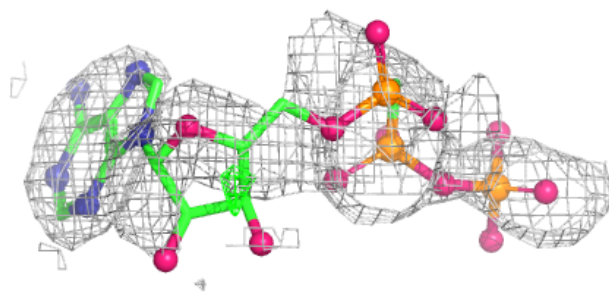
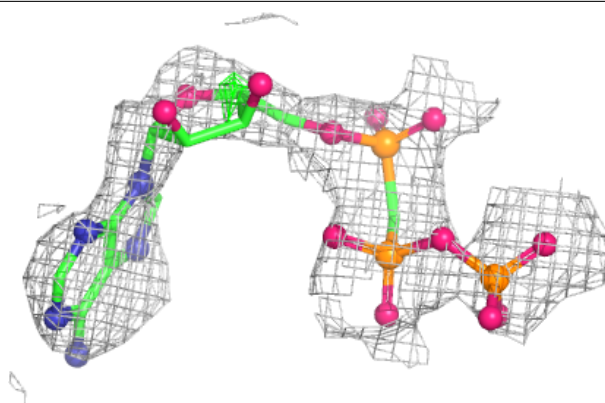
Electron density around APC G 602:

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

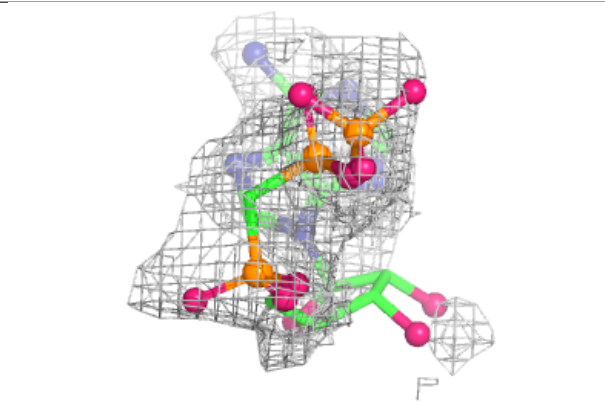
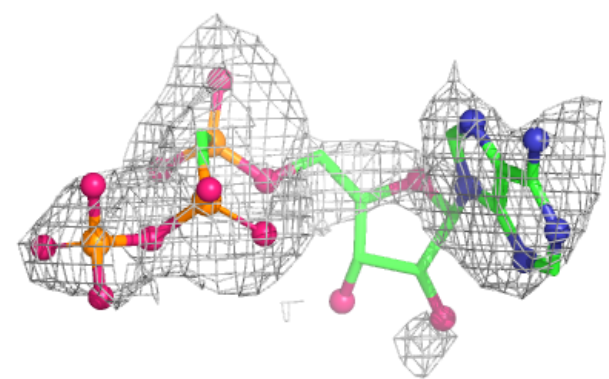
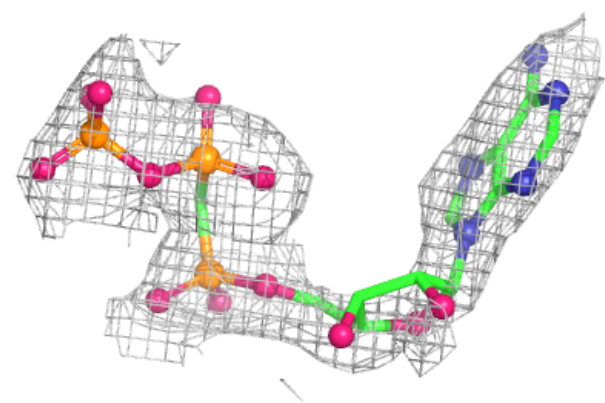


Electron density around APC H 603:

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

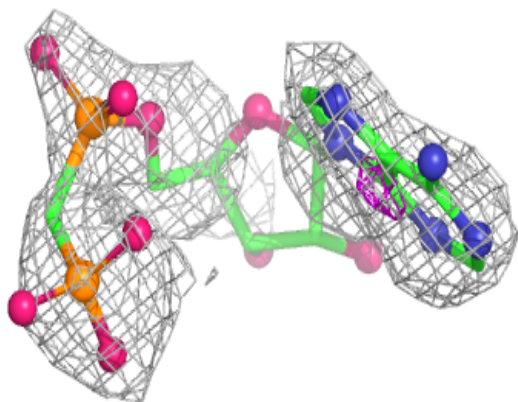
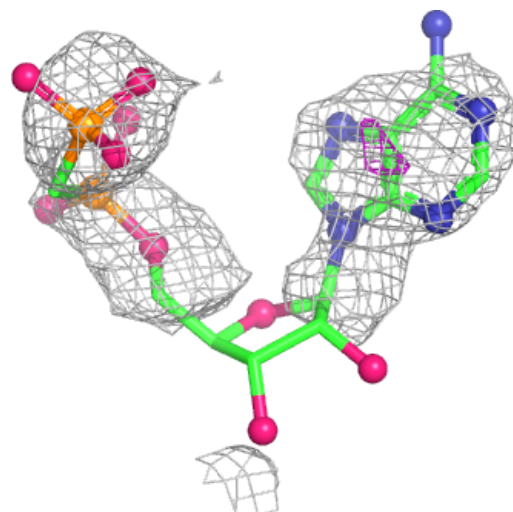
**Electron density around APC B 602:**

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



Electron density around APC A 602:

$2mF_o-DF_c$ (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.