



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:21 AM UTC

PDB ID : 6GS8 / pdb_00006gs8
Title : Crystal structure of SmbA in complex with c-di-GMP
Authors : Dubey, B.N.; Schirmer, T.
Deposited on : 2018-06-13
Resolution : 2.80 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

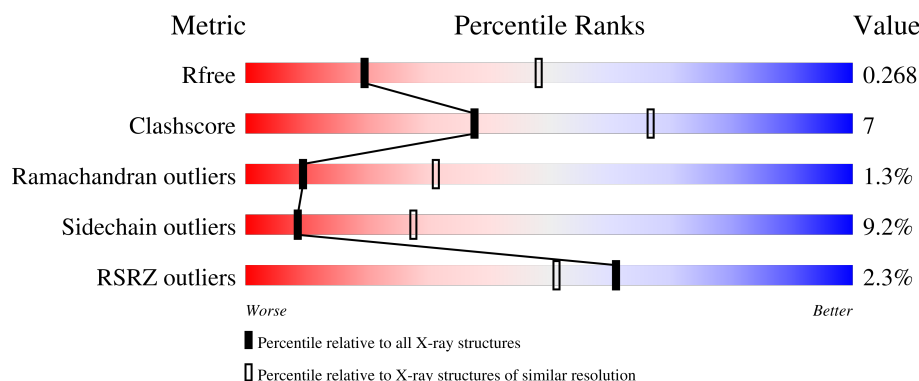
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	308	 77% 15% 5% 2% 1%
1	B	308	 75% 19% 2% 2% 1%
1	C	308	 74% 15% 8% 4% 1%
1	D	308	 74% 13% 5% 3% 1%
1	E	308	 76% 12% 8% 1% 1%

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Mol	Chain	Length	Quality of chain
1	F	308	3% 75% 16% • 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14110 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	301	Total	C	N	O	S	Se	0	0	0
			2345	1471	432	437	2	3			
1	B	301	Total	C	N	O	S	Se	0	0	0
			2345	1471	432	437	2	3			
1	C	284	Total	C	N	O	S	Se	0	0	0
			2208	1386	399	418	2	3			
1	D	284	Total	C	N	O	S	Se	0	0	0
			2208	1386	399	418	2	3			
1	E	284	Total	C	N	O	S	Se	0	0	0
			2208	1386	399	418	2	3			
1	F	284	Total	C	N	O	S	Se	0	0	0
			2208	1386	399	418	2	3			

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	initiating methionine	UNP Q9A5E6
A	198	VAL	-	insertion	UNP Q9A5E6
A	297	LYS	-	expression tag	UNP Q9A5E6
A	298	LEU	-	expression tag	UNP Q9A5E6
A	299	ALA	-	expression tag	UNP Q9A5E6
A	300	ALA	-	expression tag	UNP Q9A5E6
A	301	ALA	-	expression tag	UNP Q9A5E6
A	302	LEU	-	expression tag	UNP Q9A5E6
A	303	GLU	-	expression tag	UNP Q9A5E6
A	304	HIS	-	expression tag	UNP Q9A5E6
A	305	HIS	-	expression tag	UNP Q9A5E6
A	306	HIS	-	expression tag	UNP Q9A5E6
A	307	HIS	-	expression tag	UNP Q9A5E6
B	1	MSE	-	initiating methionine	UNP Q9A5E6
B	198	VAL	-	insertion	UNP Q9A5E6
B	297	LYS	-	expression tag	UNP Q9A5E6
B	298	LEU	-	expression tag	UNP Q9A5E6

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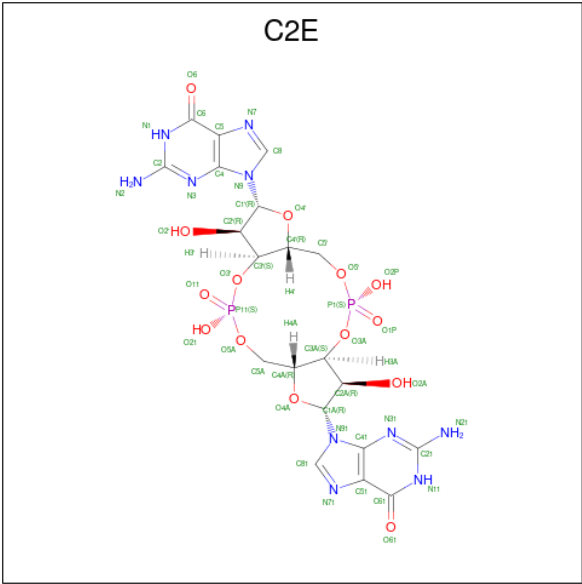
Chain	Residue	Modelled	Actual	Comment	Reference
B	299	ALA	-	expression tag	UNP Q9A5E6
B	300	ALA	-	expression tag	UNP Q9A5E6
B	301	ALA	-	expression tag	UNP Q9A5E6
B	302	LEU	-	expression tag	UNP Q9A5E6
B	303	GLU	-	expression tag	UNP Q9A5E6
B	304	HIS	-	expression tag	UNP Q9A5E6
B	305	HIS	-	expression tag	UNP Q9A5E6
B	306	HIS	-	expression tag	UNP Q9A5E6
B	307	HIS	-	expression tag	UNP Q9A5E6
C	1	MSE	-	initiating methionine	UNP Q9A5E6
C	198	VAL	-	insertion	UNP Q9A5E6
C	297	LYS	-	expression tag	UNP Q9A5E6
C	298	LEU	-	expression tag	UNP Q9A5E6
C	299	ALA	-	expression tag	UNP Q9A5E6
C	300	ALA	-	expression tag	UNP Q9A5E6
C	301	ALA	-	expression tag	UNP Q9A5E6
C	302	LEU	-	expression tag	UNP Q9A5E6
C	303	GLU	-	expression tag	UNP Q9A5E6
C	304	HIS	-	expression tag	UNP Q9A5E6
C	305	HIS	-	expression tag	UNP Q9A5E6
C	306	HIS	-	expression tag	UNP Q9A5E6
C	307	HIS	-	expression tag	UNP Q9A5E6
D	1	MSE	-	initiating methionine	UNP Q9A5E6
D	198	VAL	-	insertion	UNP Q9A5E6
D	297	LYS	-	expression tag	UNP Q9A5E6
D	298	LEU	-	expression tag	UNP Q9A5E6
D	299	ALA	-	expression tag	UNP Q9A5E6
D	300	ALA	-	expression tag	UNP Q9A5E6
D	301	ALA	-	expression tag	UNP Q9A5E6
D	302	LEU	-	expression tag	UNP Q9A5E6
D	303	GLU	-	expression tag	UNP Q9A5E6
D	304	HIS	-	expression tag	UNP Q9A5E6
D	305	HIS	-	expression tag	UNP Q9A5E6
D	306	HIS	-	expression tag	UNP Q9A5E6
D	307	HIS	-	expression tag	UNP Q9A5E6
E	1	MSE	-	initiating methionine	UNP Q9A5E6
E	198	VAL	-	insertion	UNP Q9A5E6
E	297	LYS	-	expression tag	UNP Q9A5E6
E	298	LEU	-	expression tag	UNP Q9A5E6
E	299	ALA	-	expression tag	UNP Q9A5E6
E	300	ALA	-	expression tag	UNP Q9A5E6
E	301	ALA	-	expression tag	UNP Q9A5E6

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Chain	Residue	Modelled	Actual	Comment	Reference
E	302	LEU	-	expression tag	UNP Q9A5E6
E	303	GLU	-	expression tag	UNP Q9A5E6
E	304	HIS	-	expression tag	UNP Q9A5E6
E	305	HIS	-	expression tag	UNP Q9A5E6
E	306	HIS	-	expression tag	UNP Q9A5E6
E	307	HIS	-	expression tag	UNP Q9A5E6
F	1	MSE	-	initiating methionine	UNP Q9A5E6
F	198	VAL	-	insertion	UNP Q9A5E6
F	297	LYS	-	expression tag	UNP Q9A5E6
F	298	LEU	-	expression tag	UNP Q9A5E6
F	299	ALA	-	expression tag	UNP Q9A5E6
F	300	ALA	-	expression tag	UNP Q9A5E6
F	301	ALA	-	expression tag	UNP Q9A5E6
F	302	LEU	-	expression tag	UNP Q9A5E6
F	303	GLU	-	expression tag	UNP Q9A5E6
F	304	HIS	-	expression tag	UNP Q9A5E6
F	305	HIS	-	expression tag	UNP Q9A5E6
F	306	HIS	-	expression tag	UNP Q9A5E6
F	307	HIS	-	expression tag	UNP Q9A5E6

- Molecule 2 is 9,9'-[(2R,3R,3aS,5S,7aR,9R,10R,10aS,12S,14aR)-3,5,10,12-tetrahydroxy-5,12-dioxidoctahydro-2H,7H-difuro[3,2-d:3',2'-j][1,3,7,9,2,8]tetraoxadiphosphacyclododecine-2,9-diyl]bis(2-amino-1,9-dihydro-6H-purin-6-one) (CCD ID: C2E) (formula: C₂₀H₂₄N₁₀O₁₄P₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			46	20	10	14	2		
2	A	1	Total	C	N	O	P	0	0
			46	20	10	14	2		
2	B	1	Total	C	N	O	P	0	0
			46	20	10	14	2		
2	B	1	Total	C	N	O	P	0	0
			46	20	10	14	2		
2	C	1	Total	C	N	O	P	0	0
			46	20	10	14	2		
2	C	1	Total	C	N	O	P	0	0
			46	20	10	14	2		
2	D	1	Total	C	N	O	P	0	0
			46	20	10	14	2		
2	D	1	Total	C	N	O	P	0	0
			46	20	10	14	2		
2	E	1	Total	C	N	O	P	0	0
			46	20	10	14	2		
2	E	1	Total	C	N	O	P	0	0
			46	20	10	14	2		
2	F	1	Total	C	N	O	P	46	0
			46	20	10	14	2		
2	F	1	Total	C	N	O	P	46	0
			46	20	10	14	2		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	E	1	Total	Mg	0	0
			1	1		
3	F	1	Total	Mg	0	0
			1	1		

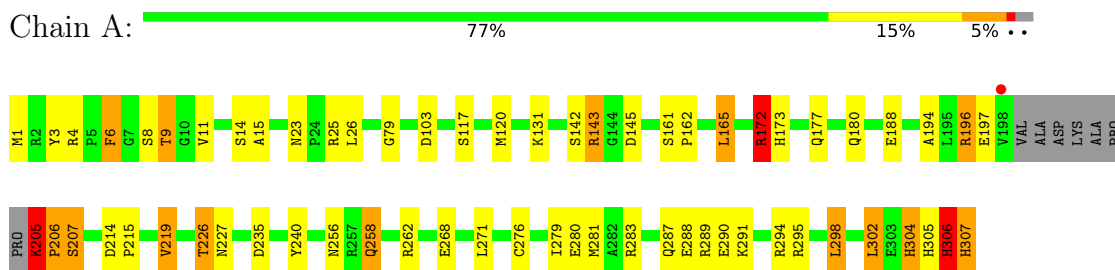
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	5	Total 5	O 5	0	0
4	B	5	Total 5	O 5	0	0
4	C	5	Total 5	O 5	0	0
4	D	5	Total 5	O 5	0	0
4	E	5	Total 5	O 5	0	0
4	F	5	Total 5	O 5	0	0

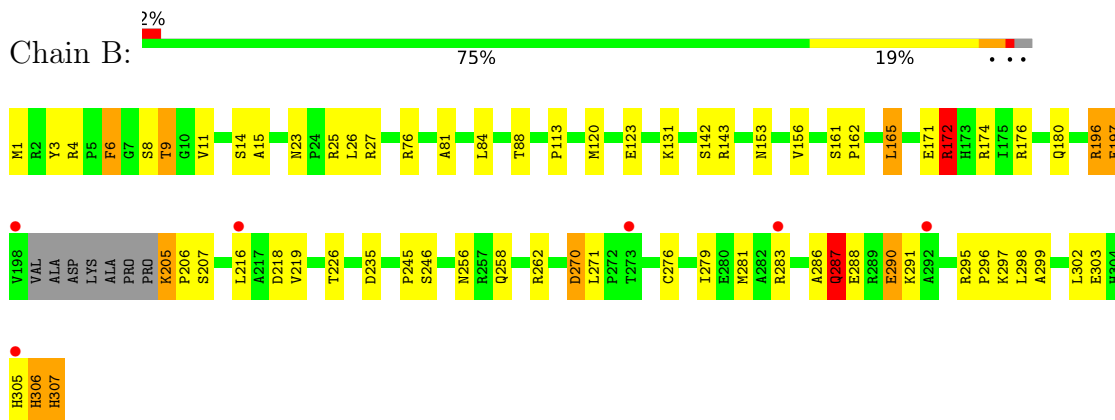
3 Residue-property plots

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.

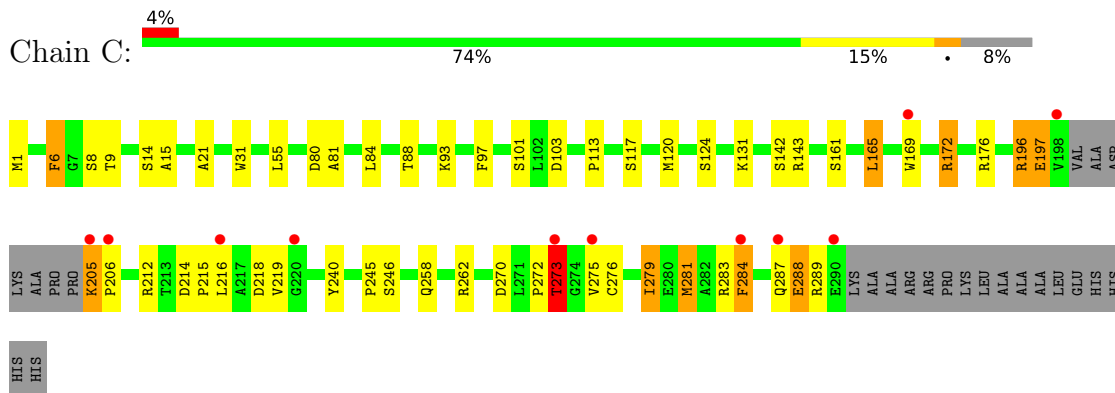
• Molecule 1: Uncharacterized protein



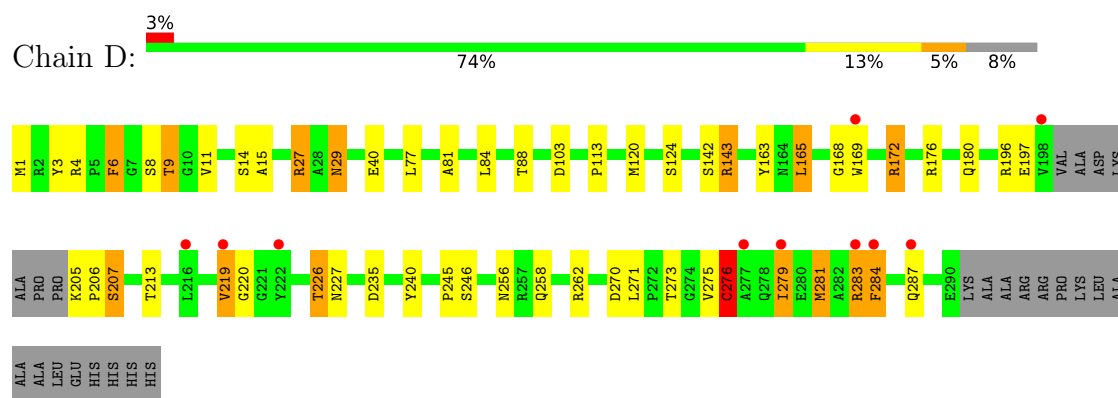
• Molecule 1: Uncharacterized protein



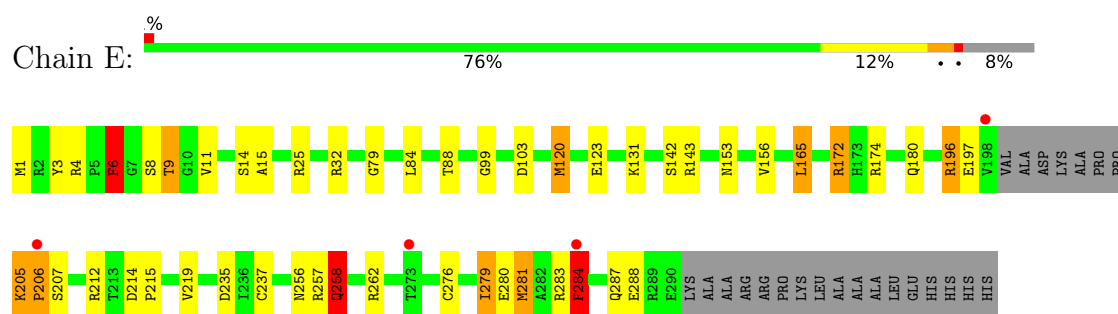
• Molecule 1: Uncharacterized protein



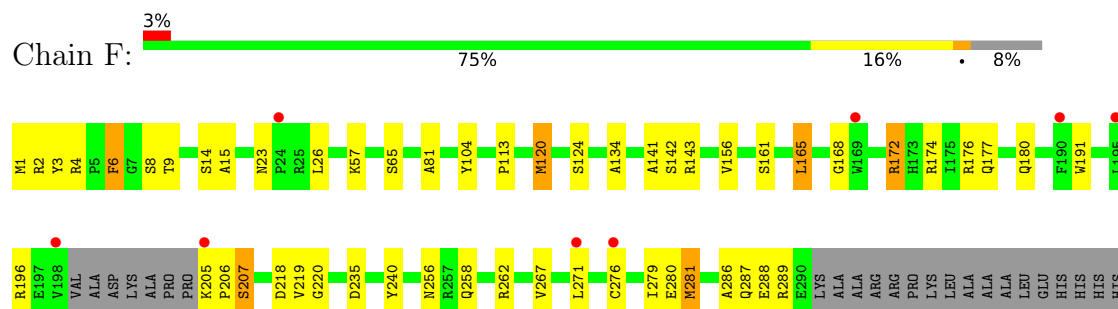
- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	127.91Å 87.42Å 130.27Å 90.00° 119.48° 90.00°	Depositor
Resolution (Å)	29.57 – 2.80 29.57 – 2.80	Depositor EDS
% Data completeness (in resolution range)	91.8 (29.57-2.80) 91.8 (29.57-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.219 , 0.268 0.221 , 0.268	Depositor DCC
R_{free} test set	1998 reflections (3.06%)	wwPDB-VP
Wilson B-factor (Å ²)	66.8	Xtriage
Anisotropy	0.464	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 36.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.002 for l,k,-h-l 0.002 for -h-l,k,h 0.014 for h,-k,-h-l 0.014 for l,-k,h 0.014 for -h-l,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14110	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.58% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, C2E

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	1.19	3/2392 (0.1%)	1.55	9/3242 (0.3%)
1	B	1.15	3/2391 (0.1%)	1.53	10/3239 (0.3%)
1	C	1.27	2/2250 (0.1%)	1.57	10/3051 (0.3%)
1	D	1.16	2/2250 (0.1%)	1.55	11/3051 (0.4%)
1	E	1.16	3/2250 (0.1%)	1.56	14/3051 (0.5%)
1	F	1.15	3/2250 (0.1%)	1.52	4/3051 (0.1%)
All	All	1.18	16/13783 (0.1%)	1.55	58/18685 (0.3%)

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	2
1	B	0	1
1	C	0	3
1	D	0	1
1	E	0	1
1	F	0	1
All	All	0	9

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	288	GLU	C-N	18.20	1.59	1.33
1	F	141	ALA	C-O	7.54	1.33	1.24
1	F	207	SER	C-O	-6.99	1.15	1.24
1	D	77	LEU	C-O	6.80	1.31	1.24
1	B	156	VAL	C-O	6.48	1.31	1.23
1	A	145	ASP	C-O	-6.06	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	173	HIS	CE1-NE2	5.96	1.38	1.32
1	E	32	ARG	NE-CZ	-5.87	1.26	1.33
1	C	273	THR	C-O	5.78	1.30	1.24
1	F	156	VAL	C-O	5.46	1.30	1.23
1	B	245	PRO	C-O	-5.44	1.17	1.24
1	E	156	VAL	C-O	5.26	1.29	1.23
1	E	99	GLY	C-O	5.22	1.29	1.23
1	D	219	VAL	N-CA	5.17	1.50	1.46
1	B	76	ARG	C-O	-5.14	1.17	1.24
1	A	304	HIS	CE1-NE2	5.11	1.37	1.32

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	281	MSE	CG-SE-CE	13.66	128.98	98.92
1	E	120	MSE	CG-SE-CE	13.54	128.71	98.92
1	E	258	GLN	CB-CG-CD	12.24	133.41	112.60
1	C	281	MSE	CG-SE-CE	8.06	116.65	98.92
1	D	27	ARG	CG-CD-NE	-8.05	94.28	112.00
1	E	258	GLN	CB-CA-C	7.86	125.76	110.67
1	E	205	LYS	CB-CA-C	7.84	125.01	110.10
1	C	6	PHE	CA-C-O	-7.51	113.73	122.37
1	C	272	PRO	CA-C-N	7.21	130.82	120.79
1	C	272	PRO	C-N-CA	7.21	130.82	120.79
1	D	143	ARG	CG-CD-NE	-7.05	96.48	112.00
1	C	120	MSE	CG-SE-CE	-6.92	83.70	98.92
1	C	284	PHE	CA-CB-CG	6.82	120.62	113.80
1	B	270	ASP	CA-CB-CG	6.75	119.35	112.60
1	B	6	PHE	CA-C-O	-6.71	114.65	122.37
1	B	205	LYS	CB-CA-C	6.59	122.62	110.10
1	E	284	PHE	CA-CB-CG	6.58	120.38	113.80
1	A	6	PHE	CA-C-O	-6.50	114.89	122.37
1	B	287	GLN	CB-CA-C	6.45	121.51	111.13
1	E	6	PHE	CA-C-O	-6.32	115.10	122.37
1	A	226	THR	CA-CB-OG1	-6.26	100.21	109.60
1	D	281	MSE	CG-SE-CE	6.24	112.66	98.92
1	D	270	ASP	CA-CB-CG	6.19	118.79	112.60
1	A	79	GLY	CA-C-O	-6.18	117.97	122.23
1	F	6	PHE	CA-C-O	-6.14	114.99	122.41
1	C	270	ASP	CA-CB-CG	6.01	118.61	112.60
1	B	143	ARG	CG-CD-NE	-6.01	98.78	112.00
1	A	172	ARG	CG-CD-NE	-5.99	98.83	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	281	MSE	CG-SE-CE	5.98	112.07	98.92
1	A	289	ARG	N-CA-C	-5.97	104.74	112.68
1	D	6	PHE	CA-C-O	-5.92	115.57	122.37
1	B	216	LEU	CA-C-O	-5.76	114.93	120.92
1	E	143	ARG	CB-CA-C	5.71	119.93	109.99
1	A	143	ARG	CB-CG-CD	5.67	124.35	111.30
1	D	143	ARG	CB-CG-CD	5.64	124.28	111.30
1	A	205	LYS	CB-CA-C	5.60	120.75	110.10
1	C	273	THR	CB-CA-C	5.50	119.67	110.92
1	C	143	ARG	CB-CA-C	5.46	119.49	109.99
1	B	226	THR	CA-CB-OG1	-5.44	101.44	109.60
1	E	237	CYS	O-C-N	5.38	128.52	122.22
1	D	284	PHE	CA-CB-CG	5.38	119.18	113.80
1	D	226	THR	CA-CB-OG1	-5.33	101.61	109.60
1	E	143	ARG	CB-CG-CD	5.32	123.53	111.30
1	D	29	ASN	N-CA-CB	5.32	118.04	110.06
1	B	290	GLU	CB-CG-CD	5.31	121.63	112.60
1	A	143	ARG	CB-CA-C	5.31	119.77	110.64
1	B	172	ARG	CG-CD-NE	-5.30	100.35	112.00
1	D	276	CYS	CB-CA-C	5.26	119.14	110.88
1	B	143	ARG	CB-CG-CD	5.24	123.36	111.30
1	E	143	ARG	CG-CD-NE	-5.22	100.52	112.00
1	F	143	ARG	CG-CD-NE	-5.21	100.55	112.00
1	D	143	ARG	N-CA-C	-5.17	105.85	113.61
1	E	280	GLU	CA-C-O	-5.16	115.38	120.70
1	A	143	ARG	CG-CD-NE	-5.13	100.72	112.00
1	E	79	GLY	CA-C-O	-5.09	117.91	121.88
1	E	172	ARG	CG-CD-NE	-5.06	100.87	112.00
1	F	120	MSE	CB-CG-SE	5.05	127.85	112.70
1	C	80	ASP	CB-CA-C	5.03	117.05	110.06

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	306	HIS	Peptide
1	A	6	PHE	Peptide
1	B	6	PHE	Peptide
1	C	288	GLU	Mainchain
1	C	289	ARG	Peptide
1	C	6	PHE	Peptide
1	D	220	GLY	Peptide

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Mol	Chain	Res	Type	Group
1	E	6	PHE	Peptide
1	F	6	PHE	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	2345	0	2313	30	0
1	B	2345	0	2312	39	0
1	C	2208	0	2173	41	0
1	D	2208	0	2173	43	0
1	E	2208	0	2173	35	0
1	F	2208	0	2173	30	0
2	A	92	0	44	1	0
2	B	92	0	44	0	0
2	C	92	0	44	1	0
2	D	92	0	44	2	0
2	E	92	0	44	2	0
2	F	92	0	44	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
4	C	5	0	0	0	0
4	D	5	0	0	0	0
4	E	5	0	0	0	0
4	F	5	0	0	0	0
All	All	14110	0	13581	192	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 (192) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:286:ALA:C	1:B:287:GLN:HA	1.79	1.07
1:D:172:ARG:NH2	1:D:240:TYR:OH	1.90	1.02
1:E:284:PHE:CZ	1:F:180:GLN:HG3	1.94	1.02
1:C:172:ARG:NH2	1:C:240:TYR:OH	1.94	1.00
1:E:256:ASN:ND2	1:E:258:GLN:HG2	1.80	0.96
1:B:286:ALA:C	1:B:287:GLN:CA	2.42	0.92
1:E:256:ASN:HD22	1:E:258:GLN:HG2	1.38	0.87
1:C:273:THR:HA	1:D:273:THR:HA	1.65	0.78
1:B:303:GLU:O	1:B:305:HIS:O	2.02	0.76
1:F:2:ARG:NH1	1:F:267:VAL:O	2.19	0.76
1:E:1:MSE:HE2	1:E:15:ALA:H	1.50	0.75
1:A:1:MSE:HE2	1:A:15:ALA:H	1.50	0.75
1:C:1:MSE:HE2	1:C:15:ALA:H	1.51	0.75
1:A:1:MSE:HE2	1:A:15:ALA:N	2.02	0.74
1:D:1:MSE:HE2	1:D:15:ALA:N	2.02	0.74
1:B:1:MSE:HE2	1:B:15:ALA:N	2.02	0.74
1:C:1:MSE:HE2	1:C:15:ALA:N	2.02	0.74
1:C:284:PHE:HD2	1:C:287:GLN:NE2	1.86	0.74
1:D:1:MSE:HE2	1:D:15:ALA:H	1.51	0.73
1:D:172:ARG:HH11	1:D:172:ARG:HG2	1.51	0.73
1:E:1:MSE:HE2	1:E:15:ALA:N	2.03	0.73
1:B:1:MSE:HE2	1:B:15:ALA:H	1.53	0.73
1:F:1:MSE:HE2	1:F:15:ALA:N	2.04	0.72
1:F:1:MSE:HE2	1:F:15:ALA:H	1.53	0.72
1:F:23:ASN:ND2	1:F:26:LEU:HD13	2.03	0.72
1:B:297:LYS:O	1:B:297:LYS:HD2	1.89	0.71
1:C:219:VAL:HG12	1:C:219:VAL:O	1.89	0.70
1:C:273:THR:CG2	1:D:273:THR:HG22	2.22	0.70
1:A:302:LEU:O	1:A:305:HIS:O	2.09	0.70
1:D:163:TYR:OH	1:D:172:ARG:NH2	2.25	0.70
1:B:286:ALA:C	1:B:287:GLN:N	2.50	0.69
1:C:245:PRO:HB3	1:D:284:PHE:CD2	2.27	0.69
1:F:220:GLY:HA3	1:F:289:ARG:HH22	1.59	0.68
1:F:218:ASP:OD2	1:F:289:ARG:NH2	2.28	0.67
1:A:180:GLN:HG2	1:B:288:GLU:HG3	1.75	0.66
1:B:286:ALA:O	1:B:287:GLN:HA	1.98	0.63
1:F:1:MSE:CE	1:F:15:ALA:HB3	2.28	0.62
1:A:1:MSE:CE	1:A:15:ALA:HB3	2.29	0.62
1:B:23:ASN:ND2	1:B:26:LEU:HD13	2.15	0.61
1:B:218:ASP:OD2	1:B:291:LYS:CE	2.48	0.61
1:D:1:MSE:CE	1:D:15:ALA:HB3	2.30	0.61
1:E:256:ASN:HD22	1:E:258:GLN:CG	2.13	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:81:ALA:HB3	1:C:113:PRO:HB3	1.84	0.60
1:B:1:MSE:CE	1:B:15:ALA:HB3	2.34	0.57
1:D:143:ARG:NH2	2:D:401:C2E:O6	2.38	0.57
1:E:257:ARG:HB2	1:E:258:GLN:NE2	2.19	0.57
1:E:1:MSE:CE	1:E:15:ALA:HB3	2.34	0.56
1:D:172:ARG:HH11	1:D:172:ARG:CG	2.19	0.56
1:B:218:ASP:OD2	1:B:291:LYS:HE2	2.07	0.55
1:C:172:ARG:NH2	1:C:240:TYR:HH	2.05	0.55
1:C:172:ARG:HG2	1:C:172:ARG:HH11	1.71	0.55
1:E:212:ARG:HG3	2:E:402:C2E:O11	2.07	0.55
1:C:172:ARG:HG2	1:C:172:ARG:NH1	2.21	0.55
1:D:172:ARG:HE	1:D:283:ARG:NH2	2.04	0.55
1:C:169:TRP:HE1	1:D:169:TRP:HE1	1.54	0.55
1:E:205:LYS:HB3	1:E:206:PRO:HD3	1.87	0.55
1:E:123:GLU:OE2	1:E:153:ASN:ND2	2.42	0.53
1:F:172:ARG:NH2	1:F:240:TYR:OH	2.42	0.52
1:B:172:ARG:HG2	1:B:172:ARG:NH1	2.23	0.52
1:F:218:ASP:OD2	1:F:286:ALA:HA	2.09	0.52
2:D:401:C2E:H8	2:D:401:C2E:O5'	2.09	0.52
1:C:1:MSE:CE	1:C:15:ALA:HB3	2.40	0.52
1:A:305:HIS:O	1:A:306:HIS:HB2	2.10	0.51
1:B:165:LEU:HD22	1:B:165:LEU:H	1.76	0.51
1:B:123:GLU:OE2	1:B:153:ASN:ND2	2.41	0.51
1:B:256:ASN:HD21	1:B:258:GLN:HG2	1.76	0.51
1:B:307:HIS:ND1	1:B:307:HIS:C	2.69	0.51
1:D:172:ARG:HE	1:D:283:ARG:HH22	1.57	0.51
1:E:257:ARG:HB2	1:E:258:GLN:HE21	1.76	0.50
1:E:284:PHE:HA	1:E:287:GLN:CD	2.36	0.50
1:D:172:ARG:HG2	1:D:172:ARG:NH1	2.25	0.50
1:D:172:ARG:NE	1:D:283:ARG:HH22	2.09	0.50
1:D:281:MSE:O	1:D:284:PHE:HB3	2.11	0.50
1:C:165:LEU:HB2	1:C:279:ILE:CG2	2.42	0.50
1:D:172:ARG:CG	1:D:172:ARG:NH1	2.73	0.50
1:C:273:THR:HB	1:D:273:THR:HA	1.94	0.50
1:D:256:ASN:HD21	1:D:258:GLN:HG2	1.77	0.49
1:E:284:PHE:CE2	1:F:180:GLN:HG3	2.44	0.49
1:E:256:ASN:HD21	1:E:258:GLN:HG2	1.74	0.49
1:F:256:ASN:ND2	1:F:258:GLN:HG2	2.27	0.48
1:A:194:ALA:HB1	1:A:219:VAL:HG11	1.95	0.48
1:E:180:GLN:HG2	1:F:288:GLU:HG2	1.96	0.48
1:E:205:LYS:HB3	1:E:206:PRO:CD	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:3:TYR:C	1:F:4:ARG:HG3	2.37	0.48
1:C:284:PHE:CD2	1:C:287:GLN:NE2	2.73	0.48
1:F:235:ASP:OD1	1:F:262:ARG:NH1	2.47	0.48
1:B:296:PRO:HB2	1:B:299:ALA:H	1.78	0.48
1:A:307:HIS:ND1	1:A:307:HIS:C	2.71	0.47
1:E:235:ASP:OD1	1:E:262:ARG:NH1	2.47	0.47
1:E:276:CYS:SG	1:F:276:CYS:HB3	2.54	0.47
1:A:240:TYR:OH	1:A:283:ARG:NH2	2.47	0.47
1:C:21:ALA:O	1:C:31:TRP:CH2	2.67	0.47
1:C:275:VAL:O	1:C:279:ILE:HG13	2.14	0.47
1:E:288:GLU:HB2	1:F:180:GLN:OE1	2.15	0.47
1:C:273:THR:HB	1:D:273:THR:HG22	1.96	0.47
1:E:284:PHE:CE2	1:F:176:ARG:HB3	2.49	0.46
1:C:165:LEU:HD12	1:C:279:ILE:HD13	1.98	0.46
1:B:235:ASP:OD1	1:B:262:ARG:NH1	2.49	0.46
1:A:172:ARG:HG2	1:A:172:ARG:NH1	2.31	0.46
1:D:6:PHE:O	1:D:9:THR:OG1	2.33	0.46
1:F:172:ARG:NH1	1:F:172:ARG:HG2	2.31	0.46
1:D:176:ARG:HG2	1:D:246:SER:HB3	1.98	0.45
1:B:81:ALA:HB3	1:B:113:PRO:HB3	1.98	0.45
1:A:271:LEU:HD23	1:A:271:LEU:HA	1.87	0.45
1:C:165:LEU:HB2	1:C:279:ILE:HG23	1.97	0.45
1:A:3:TYR:C	1:A:4:ARG:HG3	2.42	0.45
1:A:276:CYS:SG	1:B:276:CYS:HB3	2.57	0.45
1:A:287:GLN:HE21	1:B:180:GLN:HE22	1.65	0.45
1:A:305:HIS:O	1:A:306:HIS:CB	2.64	0.45
1:C:273:THR:CB	1:D:273:THR:HG22	2.45	0.45
1:C:284:PHE:CB	1:D:245:PRO:HB2	2.47	0.45
1:E:284:PHE:CD2	1:F:176:ARG:HB3	2.51	0.45
1:A:9:THR:HB	1:A:11:VAL:H	1.82	0.45
1:A:180:GLN:HG2	1:B:288:GLU:CG	2.44	0.45
1:C:84:LEU:HD22	1:C:88:THR:HG21	1.99	0.45
1:C:276:CYS:SG	1:D:276:CYS:HB3	2.57	0.45
1:B:9:THR:HB	1:B:11:VAL:H	1.82	0.44
1:C:205:LYS:HB2	1:C:205:LYS:NZ	2.32	0.44
1:A:23:ASN:ND2	1:A:26:LEU:HD13	2.32	0.44
1:F:165:LEU:HD22	1:F:165:LEU:H	1.83	0.44
1:A:165:LEU:H	1:A:165:LEU:HD22	1.81	0.44
1:B:303:GLU:O	1:B:305:HIS:C	2.61	0.44
1:F:256:ASN:HD21	1:F:258:GLN:HG2	1.82	0.44
1:B:84:LEU:HD22	1:B:88:THR:HG21	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:296:PRO:HB2	1:B:299:ALA:HB3	2.00	0.44
1:C:273:THR:CA	1:D:273:THR:HA	2.42	0.44
1:C:273:THR:HG22	1:D:273:THR:HG22	1.97	0.44
1:E:288:GLU:OE2	1:F:177:GLN:NE2	2.51	0.44
1:A:256:ASN:HD21	1:A:258:GLN:HG2	1.83	0.44
1:E:172:ARG:HG2	1:E:172:ARG:NH1	2.33	0.43
1:C:214:ASP:HB2	1:C:215:PRO:HD3	2.01	0.43
2:C:402:C2E:H5'2	2:C:402:C2E:H512	2.01	0.43
1:B:3:TYR:C	1:B:4:ARG:HG3	2.44	0.43
1:B:162:PRO:HG2	1:B:171:GLU:OE2	2.19	0.43
1:E:3:TYR:C	1:E:4:ARG:HG3	2.43	0.43
1:B:256:ASN:ND2	1:B:258:GLN:HG2	2.34	0.43
1:E:6:PHE:O	1:E:9:THR:OG1	2.34	0.43
1:E:9:THR:HB	1:E:11:VAL:H	1.84	0.43
2:E:401:C2E:H81	2:E:401:C2E:O5A	2.19	0.43
1:A:205:LYS:HB3	1:A:206:PRO:CD	2.49	0.43
1:B:305:HIS:O	1:B:306:HIS:CB	2.67	0.43
1:C:93:LYS:HE2	1:C:97:PHE:CD2	2.53	0.43
1:A:288:GLU:HG2	1:B:180:GLN:HG2	2.01	0.42
1:D:226:THR:HG22	1:D:227:ASN:ND2	2.34	0.42
1:D:84:LEU:HD22	1:D:88:THR:HG21	2.02	0.42
1:A:143:ARG:NH2	2:A:401:C2E:O6	2.52	0.42
1:D:9:THR:HB	1:D:11:VAL:H	1.85	0.42
1:D:284:PHE:CD2	1:D:284:PHE:O	2.73	0.42
1:C:287:GLN:HB2	1:D:180:GLN:CD	2.45	0.42
1:F:81:ALA:HB3	1:F:113:PRO:HB3	2.01	0.42
1:A:214:ASP:HB2	1:A:215:PRO:HD3	2.01	0.42
1:F:271:LEU:HD23	1:F:271:LEU:HA	1.91	0.42
1:D:235:ASP:OD1	1:D:262:ARG:NH1	2.52	0.42
1:D:271:LEU:HD23	1:D:271:LEU:HA	1.89	0.42
1:E:214:ASP:HB2	1:E:215:PRO:HD3	2.01	0.42
1:D:81:ALA:HB3	1:D:113:PRO:HB3	2.02	0.42
1:A:1:MSE:HE1	1:A:15:ALA:HB3	2.00	0.41
1:C:273:THR:HG22	1:D:273:THR:CG2	2.51	0.41
1:D:3:TYR:C	1:D:4:ARG:HG3	2.44	0.41
1:E:165:LEU:C	1:E:165:LEU:HD23	2.45	0.41
1:C:165:LEU:HD23	1:C:165:LEU:C	2.45	0.41
1:D:165:LEU:C	1:D:165:LEU:HD23	2.45	0.41
1:E:196:ARG:C	1:E:197:GLU:HG3	2.46	0.41
1:E:165:LEU:HD12	1:E:279:ILE:HD13	2.02	0.41
1:C:31:TRP:CD1	1:C:55:LEU:HA	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:172:ARG:HH21	1:F:240:TYR:HH	1.63	0.41
1:A:235:ASP:OD1	1:A:262:ARG:NH1	2.53	0.41
1:C:172:ARG:HH21	1:C:283:ARG:NH2	2.18	0.41
1:C:284:PHE:HA	1:C:287:GLN:CD	2.46	0.41
1:D:1:MSE:HE1	1:D:15:ALA:HB3	1.99	0.41
1:A:196:ARG:O	1:A:197:GLU:HG3	2.20	0.41
1:B:176:ARG:HG2	1:B:246:SER:HB3	2.02	0.41
1:E:256:ASN:HD22	1:E:258:GLN:CD	2.28	0.41
1:A:226:THR:HG22	1:A:227:ASN:ND2	2.36	0.41
1:B:172:ARG:HG2	1:B:172:ARG:HH11	1.84	0.41
1:E:180:GLN:HE22	1:F:287:GLN:HG3	1.86	0.41
1:E:284:PHE:CZ	1:F:176:ARG:O	2.74	0.41
1:C:176:ARG:HG2	1:C:246:SER:HB3	2.03	0.41
1:C:219:VAL:O	1:C:219:VAL:CG1	2.54	0.40
1:C:284:PHE:CG	1:D:245:PRO:HB2	2.55	0.40
1:F:104:TYR:HA	1:F:134:ALA:O	2.21	0.40
1:C:196:ARG:C	1:C:197:GLU:HG3	2.46	0.40
1:F:1:MSE:HE1	1:F:15:ALA:HB3	2.02	0.40
1:A:298:LEU:HD23	1:A:298:LEU:HA	1.96	0.40
1:B:305:HIS:O	1:B:306:HIS:CG	2.74	0.40
1:D:275:VAL:O	1:D:279:ILE:HG13	2.20	0.40
1:A:162:PRO:HA	1:A:188:GLU:O	2.22	0.40
1:B:23:ASN:ND2	1:B:26:LEU:CD1	2.84	0.40
1:B:196:ARG:C	1:B:197:GLU:HG3	2.47	0.40
1:B:271:LEU:HD23	1:B:271:LEU:HA	1.87	0.40
1:D:27:ARG:NH1	1:D:27:ARG:CG	2.82	0.40
1:E:84:LEU:HD22	1:E:88:THR:HG21	2.04	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	297/308 (96%)	273 (92%)	20 (7%)	4 (1%)	9	31
1	B	295/308 (96%)	264 (90%)	26 (9%)	5 (2%)	7	25
1	C	280/308 (91%)	255 (91%)	22 (8%)	3 (1%)	11	36
1	D	280/308 (91%)	260 (93%)	17 (6%)	3 (1%)	11	36
1	E	280/308 (91%)	261 (93%)	16 (6%)	3 (1%)	11	36
1	F	280/308 (91%)	260 (93%)	16 (6%)	4 (1%)	9	30
All	All	1712/1848 (93%)	1573 (92%)	117 (7%)	22 (1%)	9	31

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	306	HIS
1	B	295	ARG
1	B	306	HIS
1	A	206	PRO
1	B	8	SER
1	B	206	PRO
1	C	8	SER
1	C	206	PRO
1	C	216	LEU
1	D	206	PRO
1	D	207	SER
1	E	206	PRO
1	F	8	SER
1	F	168	GLY
1	F	206	PRO
1	A	8	SER
1	A	207	SER
1	B	207	SER
1	E	8	SER
1	E	207	SER
1	F	207	SER
1	D	168	GLY

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	241/243 (99%)	212 (88%)	29 (12%)	5	17
1	B	241/243 (99%)	217 (90%)	24 (10%)	7	24
1	C	229/243 (94%)	208 (91%)	21 (9%)	8	27
1	D	229/243 (94%)	208 (91%)	21 (9%)	8	27
1	E	229/243 (94%)	213 (93%)	16 (7%)	14	40
1	F	229/243 (94%)	211 (92%)	18 (8%)	11	35
All	All	1398/1458 (96%)	1269 (91%)	129 (9%)	8	27

All (129) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	THR
1	A	14	SER
1	A	25	ARG
1	A	103	ASP
1	A	117	SER
1	A	120	MSE
1	A	131	LYS
1	A	142	SER
1	A	161	SER
1	A	165	LEU
1	A	172	ARG
1	A	177	GLN
1	A	196	ARG
1	A	205	LYS
1	A	207	SER
1	A	219	VAL
1	A	258	GLN
1	A	268	GLU
1	A	279	ILE
1	A	280	GLU
1	A	281	MSE
1	A	290	GLU
1	A	291	LYS
1	A	294	ARG
1	A	295	ARG
1	A	298	LEU
1	A	302	LEU
1	A	304	HIS

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Mol	Chain	Res	Type
1	A	307	HIS
1	B	9	THR
1	B	14	SER
1	B	25	ARG
1	B	27	ARG
1	B	120	MSE
1	B	131	LYS
1	B	142	SER
1	B	161	SER
1	B	165	LEU
1	B	172	ARG
1	B	174	ARG
1	B	196	ARG
1	B	197	GLU
1	B	205	LYS
1	B	219	VAL
1	B	270	ASP
1	B	279	ILE
1	B	281	MSE
1	B	283	ARG
1	B	287	GLN
1	B	290	GLU
1	B	298	LEU
1	B	302	LEU
1	B	307	HIS
1	C	9	THR
1	C	14	SER
1	C	101	SER
1	C	103	ASP
1	C	117	SER
1	C	124	SER
1	C	131	LYS
1	C	142	SER
1	C	161	SER
1	C	165	LEU
1	C	172	ARG
1	C	196	ARG
1	C	197	GLU
1	C	205	LYS
1	C	212	ARG
1	C	218	ASP
1	C	258	GLN

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Mol	Chain	Res	Type
1	C	262	ARG
1	C	273	THR
1	C	279	ILE
1	C	281	MSE
1	D	8	SER
1	D	9	THR
1	D	14	SER
1	D	29	ASN
1	D	40	GLU
1	D	103	ASP
1	D	120	MSE
1	D	124	SER
1	D	142	SER
1	D	165	LEU
1	D	172	ARG
1	D	196	ARG
1	D	197	GLU
1	D	205	LYS
1	D	207	SER
1	D	213	THR
1	D	219	VAL
1	D	276	CYS
1	D	279	ILE
1	D	283	ARG
1	D	287	GLN
1	E	9	THR
1	E	14	SER
1	E	25	ARG
1	E	103	ASP
1	E	120	MSE
1	E	131	LYS
1	E	142	SER
1	E	165	LEU
1	E	174	ARG
1	E	196	ARG
1	E	219	VAL
1	E	258	GLN
1	E	279	ILE
1	E	281	MSE
1	E	283	ARG
1	E	284	PHE
1	F	9	THR

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Mol	Chain	Res	Type
1	F	14	SER
1	F	57	LYS
1	F	65	SER
1	F	120	MSE
1	F	124	SER
1	F	142	SER
1	F	161	SER
1	F	165	LEU
1	F	172	ARG
1	F	174	ARG
1	F	191	TRP
1	F	196	ARG
1	F	205	LYS
1	F	219	VAL
1	F	279	ILE
1	F	280	GLU
1	F	281	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	181	ASN
1	A	182	ASN
1	A	210	GLN
1	A	227	ASN
1	A	306	HIS
1	B	69	HIS
1	B	114	GLN
1	B	180	GLN
1	B	181	ASN
1	B	182	ASN
1	B	193	GLN
1	B	210	GLN
1	B	306	HIS
1	C	182	ASN
1	C	210	GLN
1	C	227	ASN
1	D	114	GLN
1	D	180	GLN
1	D	182	ASN
1	D	193	GLN
1	D	227	ASN

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Mol	Chain	Res	Type
1	E	177	GLN
1	E	180	GLN
1	E	181	ASN
1	E	182	ASN
1	E	193	GLN
1	E	256	ASN
1	E	258	GLN
1	F	177	GLN
1	F	182	ASN
1	F	193	GLN

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 18 ligands modelled in this entry, 6 are monoatomic - leaving 12 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	C2E	E	401	-	52,52,52	0.53	0	78,82,82	0.75	2 (2%)
2	C2E	C	401	-	52,52,52	0.68	0	78,82,82	1.06	7 (8%)
2	C2E	B	401	-	52,52,52	0.65	0	78,82,82	0.99	5 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	C2E	A	402	-	52,52,52	0.58	0	78,82,82	1.00	5 (6%)
2	C2E	A	401	-	52,52,52	0.54	0	78,82,82	0.86	2 (2%)
2	C2E	E	402	-	52,52,52	0.49	0	78,82,82	0.82	2 (2%)
2	C2E	C	402	-	52,52,52	0.79	2 (3%)	78,82,82	0.98	5 (6%)
2	C2E	D	402	-	52,52,52	0.96	3 (5%)	78,82,82	1.10	6 (7%)
2	C2E	F	402	-	52,52,52	0.94	3 (5%)	78,82,82	1.15	8 (10%)
2	C2E	F	401	-	52,52,52	1.28	7 (13%)	78,82,82	1.41	13 (16%)
2	C2E	D	401	-	52,52,52	0.80	1 (1%)	78,82,82	1.03	4 (5%)
2	C2E	B	402	-	52,52,52	0.57	0	78,82,82	0.93	3 (3%)

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	C2E	E	401	-	-	0/30/62/62	0/6/7/7
2	C2E	C	401	-	-	0/30/62/62	0/6/7/7
2	C2E	B	401	-	-	1/30/62/62	0/6/7/7
2	C2E	A	402	-	-	0/30/62/62	0/6/7/7
2	C2E	A	401	-	-	2/30/62/62	0/6/7/7
2	C2E	E	402	-	-	4/30/62/62	0/6/7/7
2	C2E	C	402	-	-	5/30/62/62	0/6/7/7
2	C2E	D	402	-	-	0/30/62/62	0/6/7/7
2	C2E	F	402	-	-	6/30/62/62	0/6/7/7
2	C2E	F	401	-	-	3/30/62/62	0/6/7/7
2	C2E	D	401	-	-	0/30/62/62	0/6/7/7
2	C2E	B	402	-	-	7/30/62/62	0/6/7/7

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	401	C2E	C1A-N91	3.92	1.58	1.47
2	F	401	C2E	O61-C61	3.09	1.29	1.23
2	F	401	C2E	C21-N11	3.05	1.45	1.37
2	F	402	C2E	C1A-N91	-2.96	1.39	1.47
2	D	401	C2E	C4-N3	2.91	1.40	1.34
2	D	402	C2E	C81-N71	2.90	1.40	1.32
2	F	402	C2E	C1'-N9	2.67	1.55	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	402	C2E	C1A-N91	2.43	1.54	1.47
2	F	402	C2E	C41-N91	-2.33	1.32	1.38
2	F	401	C2E	C4-N3	2.29	1.39	1.34
2	C	402	C2E	C81-N71	-2.22	1.26	1.32
2	F	401	C2E	P11-O3'	-2.18	1.52	1.59
2	F	401	C2E	C1'-N9	2.14	1.53	1.47
2	F	401	C2E	C2'-C3'	2.08	1.57	1.53
2	C	402	C2E	C51-C41	-2.06	1.32	1.38
2	D	402	C2E	C51-C41	2.03	1.44	1.38

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	401	C2E	O3A-P1-O1P	-4.29	96.11	109.81
2	A	402	C2E	O2P-P1-O1P	4.02	131.14	112.44
2	F	401	C2E	C61-C51-C41	3.79	124.53	118.83
2	B	402	C2E	O2A-C2A-C3A	-3.51	101.35	111.19
2	A	402	C2E	O3A-P1-O1P	-3.39	98.97	109.81
2	F	402	C2E	O5A-P11-O11	-3.20	96.24	108.94
2	F	402	C2E	C3'-C2'-C1'	3.14	106.79	99.89
2	E	401	C2E	O21-P11-O5A	3.14	121.78	107.57
2	D	401	C2E	C2A-C3A-C4A	-3.13	97.76	103.24
2	F	401	C2E	O3'-C3'-C2'	2.98	122.38	111.68
2	F	401	C2E	O2P-P1-O1P	2.97	126.26	112.44
2	A	402	C2E	O5A-P11-O11	-2.95	97.22	108.94
2	A	402	C2E	O2'-C2'-C3'	2.93	119.40	111.19
2	C	401	C2E	O3'-P11-O11	-2.82	100.80	109.81
2	B	401	C2E	O2P-P1-O3A	2.77	117.99	106.70
2	D	401	C2E	O3A-P1-O1P	-2.72	101.11	109.81
2	D	402	C2E	O2P-P1-O3A	2.68	117.61	106.70
2	A	402	C2E	O21-P11-O11	2.67	124.85	112.44
2	F	401	C2E	C2'-C1'-N9	-2.65	105.86	113.25
2	C	401	C2E	N2-C2-N1	-2.64	111.19	116.76
2	C	402	C2E	O2'-C2'-C3'	-2.63	103.83	111.19
2	C	402	C2E	O5A-P11-O11	-2.58	98.71	108.94
2	C	401	C2E	N2-C2-N3	2.58	124.70	119.67
2	D	401	C2E	O6-C6-N1	-2.56	115.29	120.11
2	D	402	C2E	O4A-C1A-N91	-2.55	102.58	108.36
2	D	401	C2E	O2A-C2A-C3A	-2.54	104.08	111.19
2	C	401	C2E	O2'-C2'-C3'	-2.52	104.14	111.19
2	B	402	C2E	O3'-P11-O11	2.49	117.78	109.81
2	F	401	C2E	C3'-C2'-C1'	2.48	105.34	99.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	401	C2E	C61-C51-N71	-2.45	125.83	130.29
2	F	401	C2E	O21-P11-O11	2.45	123.82	112.44
2	C	401	C2E	O2A-C2A-C3A	-2.44	104.35	111.19
2	F	402	C2E	N91-C41-N31	-2.44	121.08	125.95
2	C	401	C2E	O21-P11-O11	2.41	123.66	112.44
2	C	402	C2E	O3A-P1-O1P	-2.40	102.12	109.81
2	D	402	C2E	C2A-C3A-C4A	-2.37	99.09	103.24
2	A	401	C2E	O2P-P1-O3A	-2.32	97.24	106.70
2	F	402	C2E	O3'-C3'-C2'	2.32	120.00	111.68
2	F	402	C2E	O21-P11-O11	2.29	123.11	112.44
2	F	401	C2E	C2'-C3'-C4'	-2.29	99.23	103.24
2	B	402	C2E	O2P-P1-O1P	2.28	123.03	112.44
2	F	401	C2E	C51-C61-N11	-2.27	107.47	113.25
2	F	401	C2E	O3'-C3'-C4'	-2.25	102.11	110.03
2	E	402	C2E	O21-P11-O11	2.23	122.81	112.44
2	D	402	C2E	O2A-C2A-C1A	2.23	117.77	110.10
2	F	402	C2E	C51-C41-N91	2.17	109.53	105.66
2	F	402	C2E	O5'-P1-O1P	2.14	117.43	108.94
2	F	402	C2E	O2P-P1-O1P	2.12	122.32	112.44
2	A	401	C2E	P1-O5'-C5'	-2.11	109.23	121.35
2	C	402	C2E	O21-P11-O11	2.11	122.25	112.44
2	F	401	C2E	C21-N31-C41	-2.10	108.67	112.30
2	D	402	C2E	O2A-C2A-C3A	-2.08	105.36	111.19
2	C	402	C2E	O2P-P1-O1P	2.08	122.11	112.44
2	C	401	C2E	O6-C6-C5	2.06	131.97	126.53
2	D	402	C2E	O3A-P1-O1P	-2.04	103.27	109.81
2	F	401	C2E	O3A-C3A-C4A	-2.04	102.84	110.03
2	B	401	C2E	C2A-C3A-C4A	-2.02	99.70	103.24
2	E	402	C2E	O21-P11-O3'	2.02	114.91	106.70
2	E	401	C2E	O5A-P11-O11	-2.01	100.97	108.94
2	B	401	C2E	O3'-C3'-C4'	-2.00	102.97	110.03
2	B	401	C2E	O3A-C3A-C4A	-2.00	102.97	110.03
2	B	401	C2E	O3'-C3'-C2'	2.00	118.85	111.68

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	402	C2E	C5'-O5'-P1-O1P
2	B	402	C2E	C5'-O5'-P1-O3A
2	C	402	C2E	C5'-O5'-P1-O1P
2	C	402	C2E	O4'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
2	E	402	C2E	C2'-C3'-O3'-P11
2	F	402	C2E	C4'-C3'-O3'-P11
2	F	402	C2E	C2'-C3'-O3'-P11
2	E	402	C2E	C4'-C3'-O3'-P11
2	B	402	C2E	O4'-C4'-C5'-O5'
2	B	402	C2E	C3'-C4'-C5'-O5'
2	A	401	C2E	C4'-C3'-O3'-P11
2	F	401	C2E	C4A-C3A-O3A-P1
2	F	402	C2E	C4A-C3A-O3A-P1
2	C	402	C2E	C3'-C4'-C5'-O5'
2	E	402	C2E	C2A-C3A-O3A-P1
2	B	402	C2E	C5'-O5'-P1-O2P
2	C	402	C2E	C5'-O5'-P1-O2P
2	C	402	C2E	C5'-O5'-P1-O3A
2	F	402	C2E	C5A-O5A-P11-O11
2	A	401	C2E	C2'-C3'-O3'-P11
2	B	401	C2E	C4A-C3A-O3A-P1
2	F	402	C2E	C2A-C3A-O3A-P1
2	F	401	C2E	C2A-C3A-O3A-P1
2	F	402	C2E	O4A-C4A-C5A-O5A
2	B	402	C2E	C2'-C3'-O3'-P11
2	E	402	C2E	C4A-C3A-O3A-P1
2	F	401	C2E	C3A-O3A-P1-O2P
2	B	402	C2E	C4'-C3'-O3'-P11

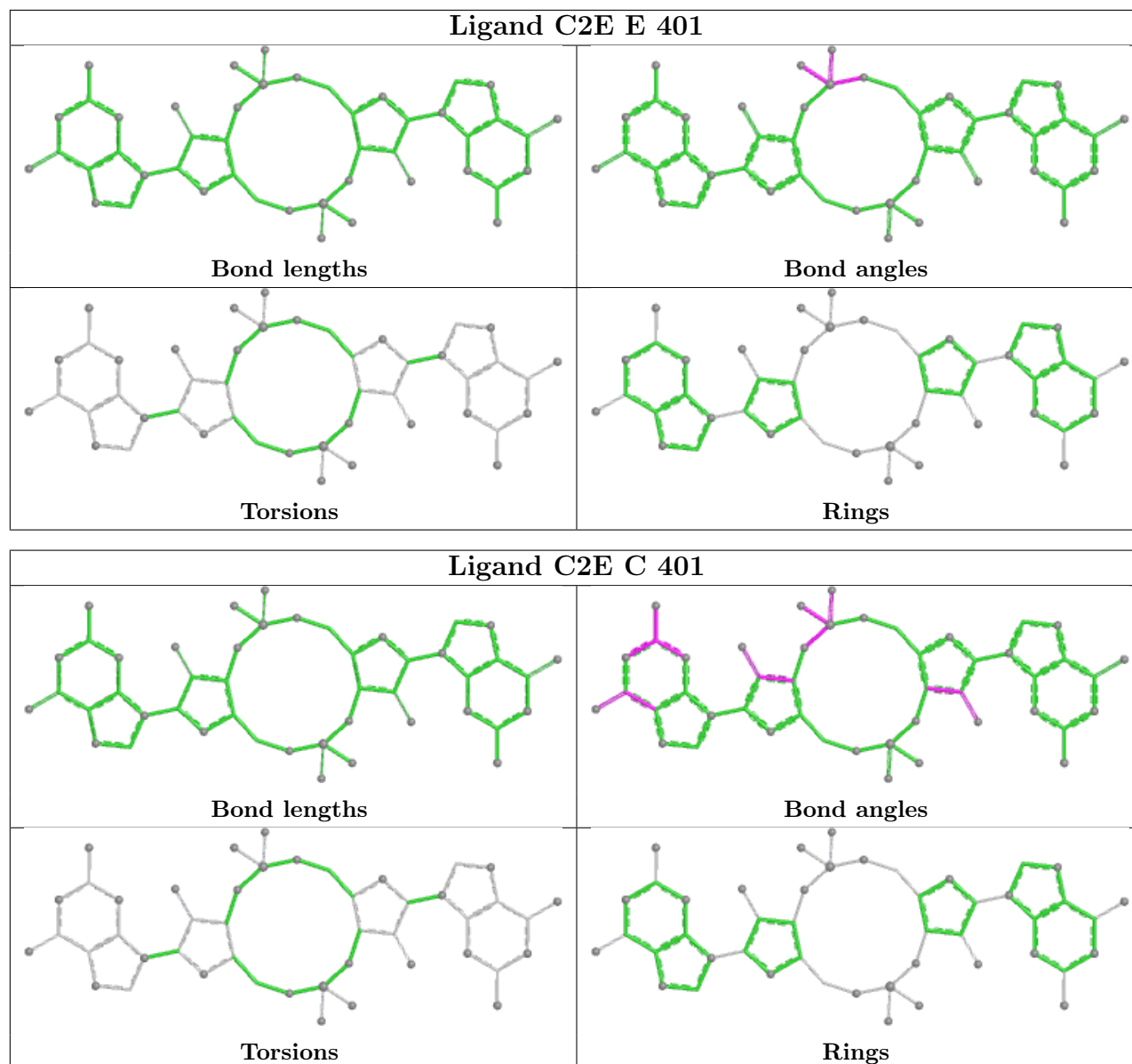
There are no ring outliers.

5 monomers are involved in 6 short contacts:

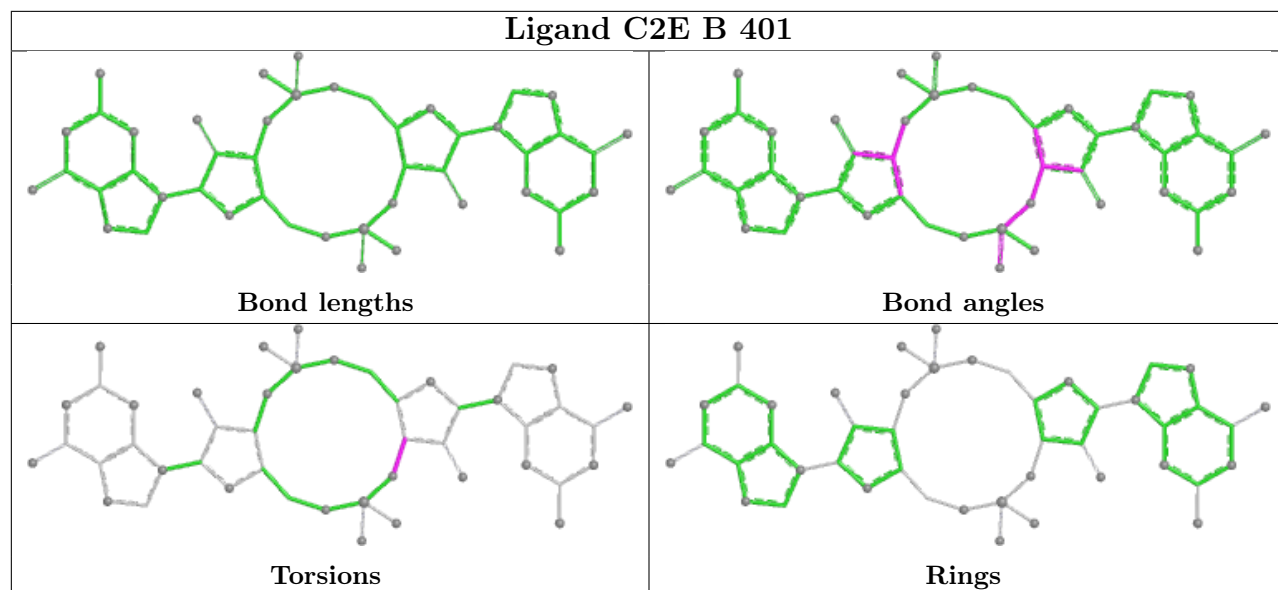
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	401	C2E	1	0
2	A	401	C2E	1	0
2	E	402	C2E	1	0
2	C	402	C2E	1	0
2	D	401	C2E	2	0

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

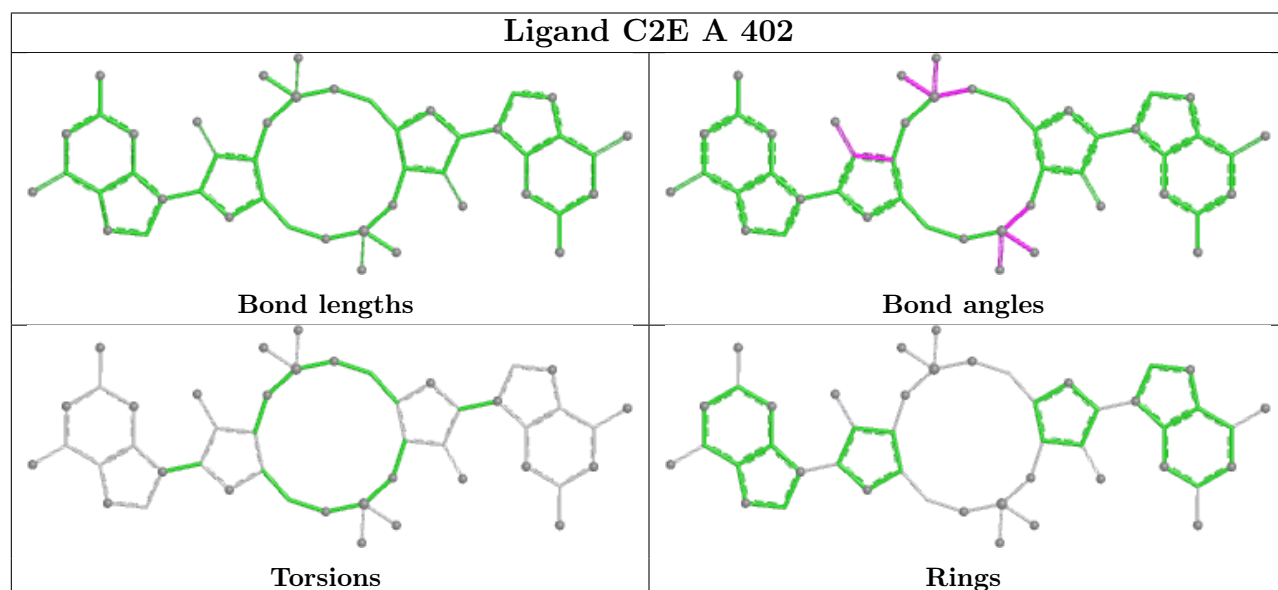
in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



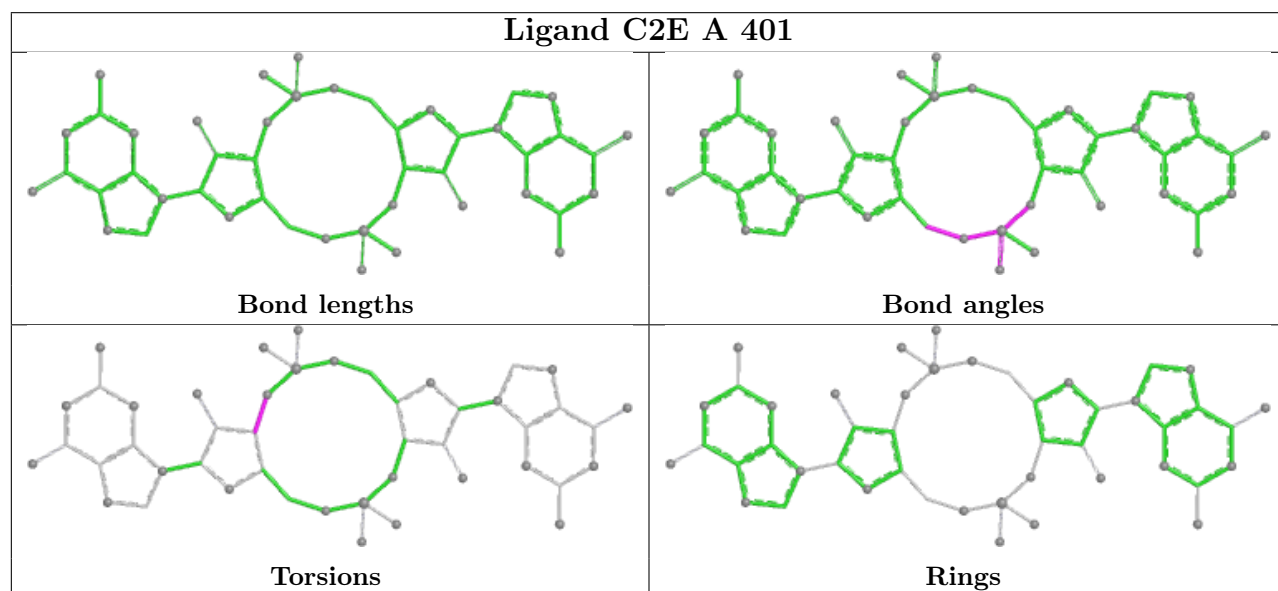
Ligand C2E B 401



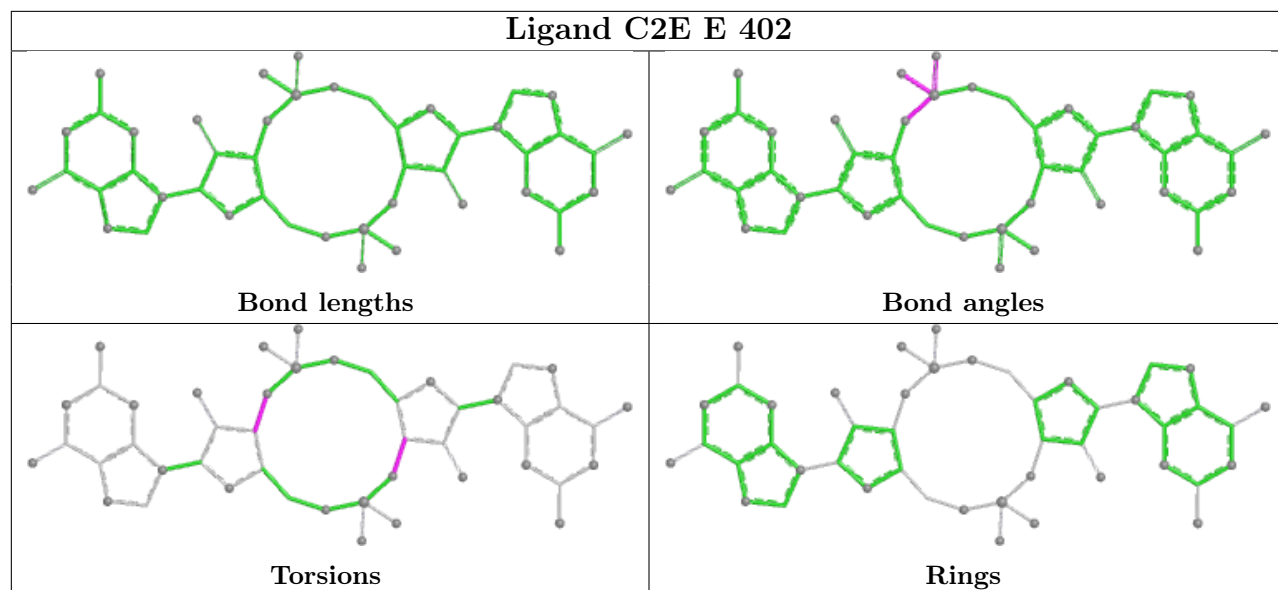
Ligand C2E A 402



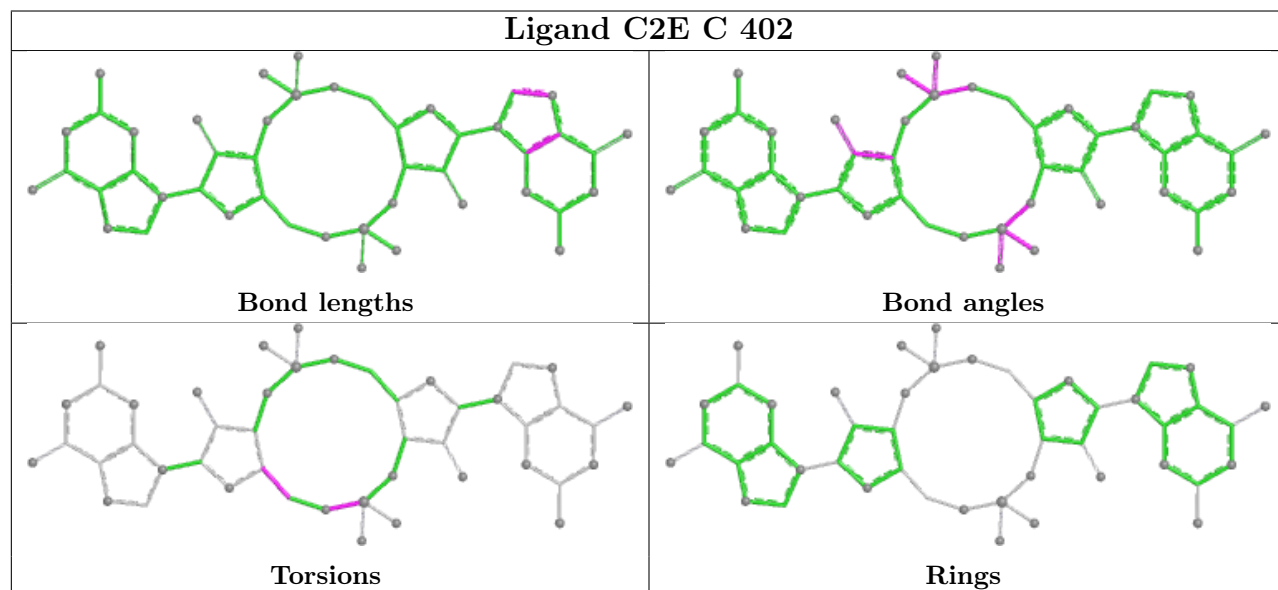
Ligand C2E A 401



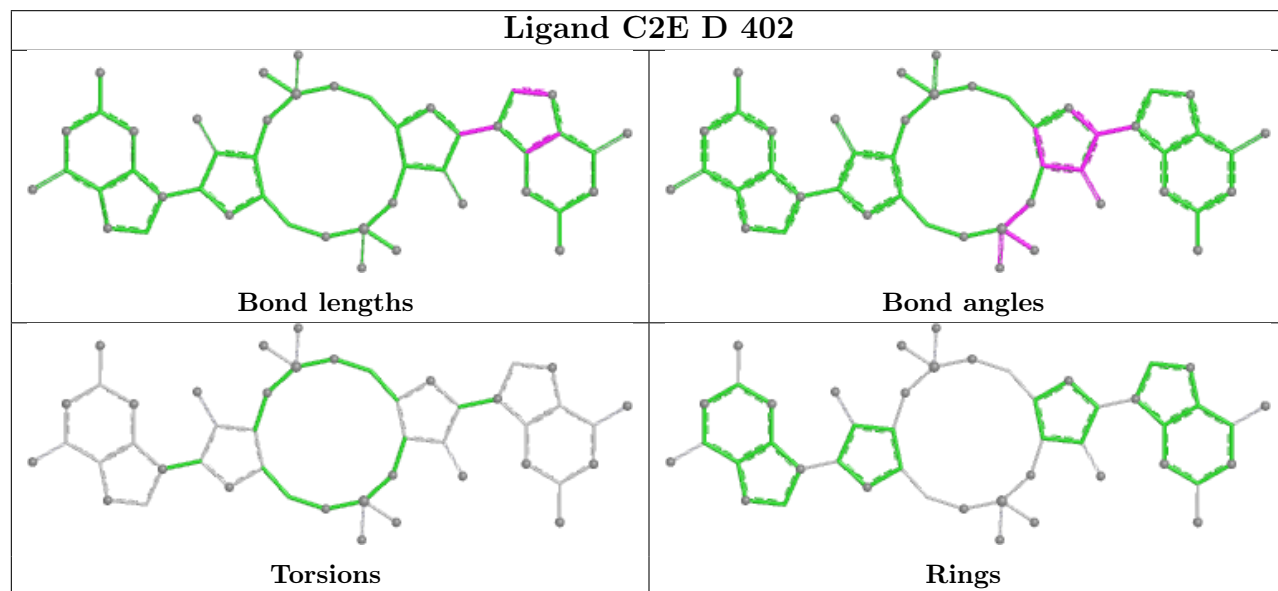
Ligand C2E E 402

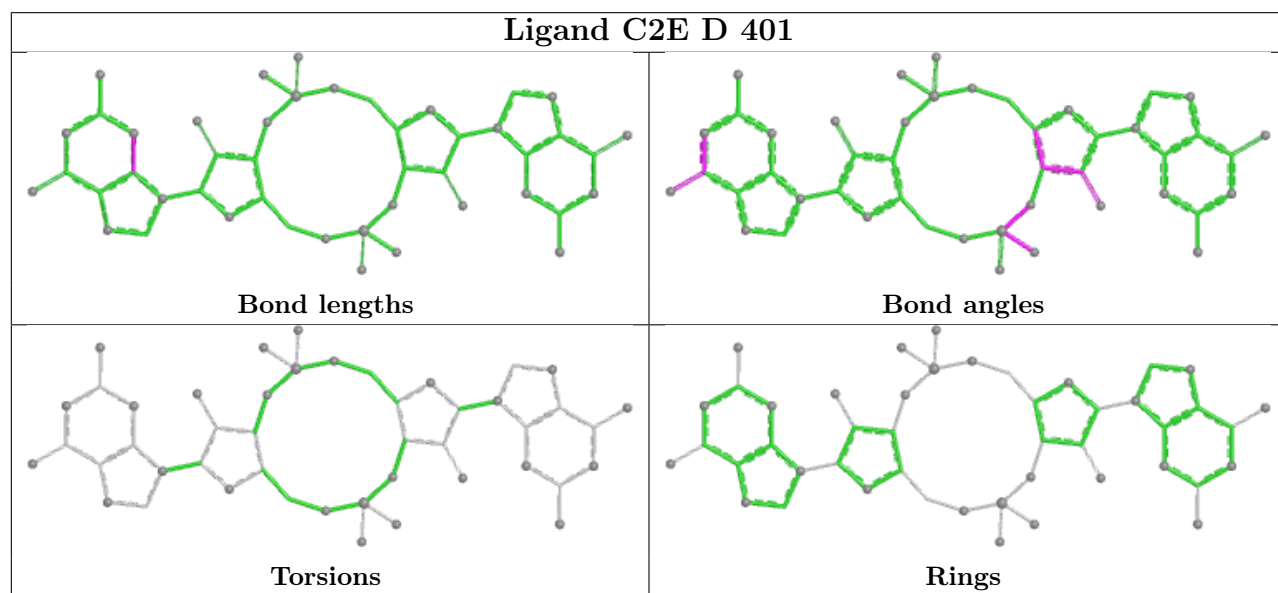
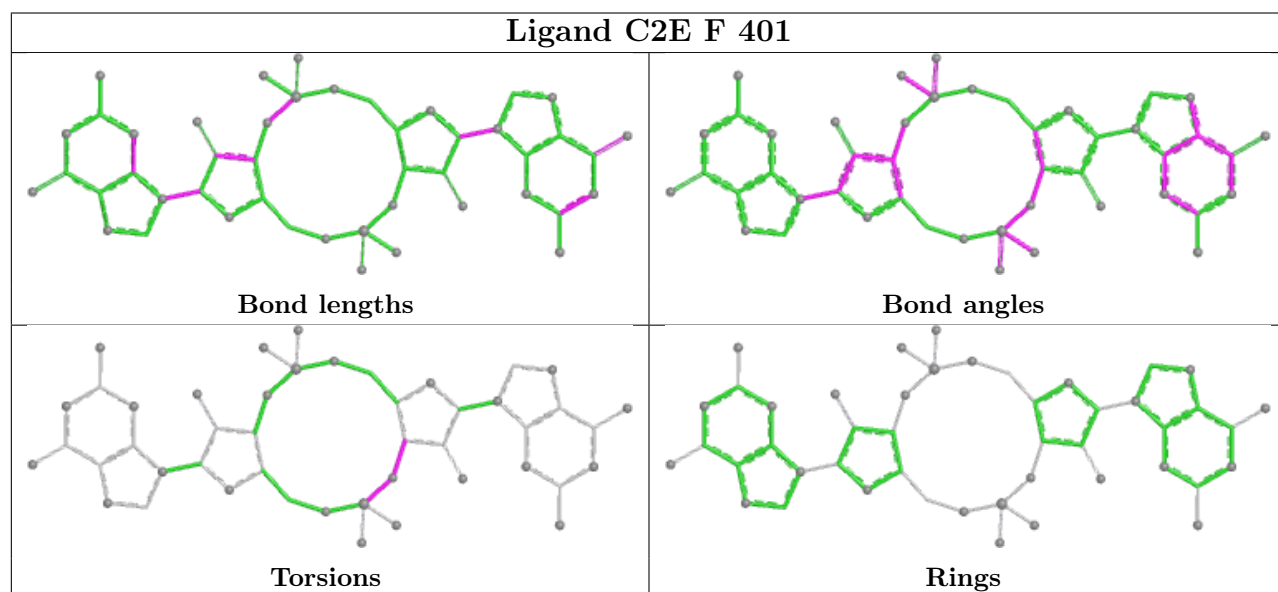
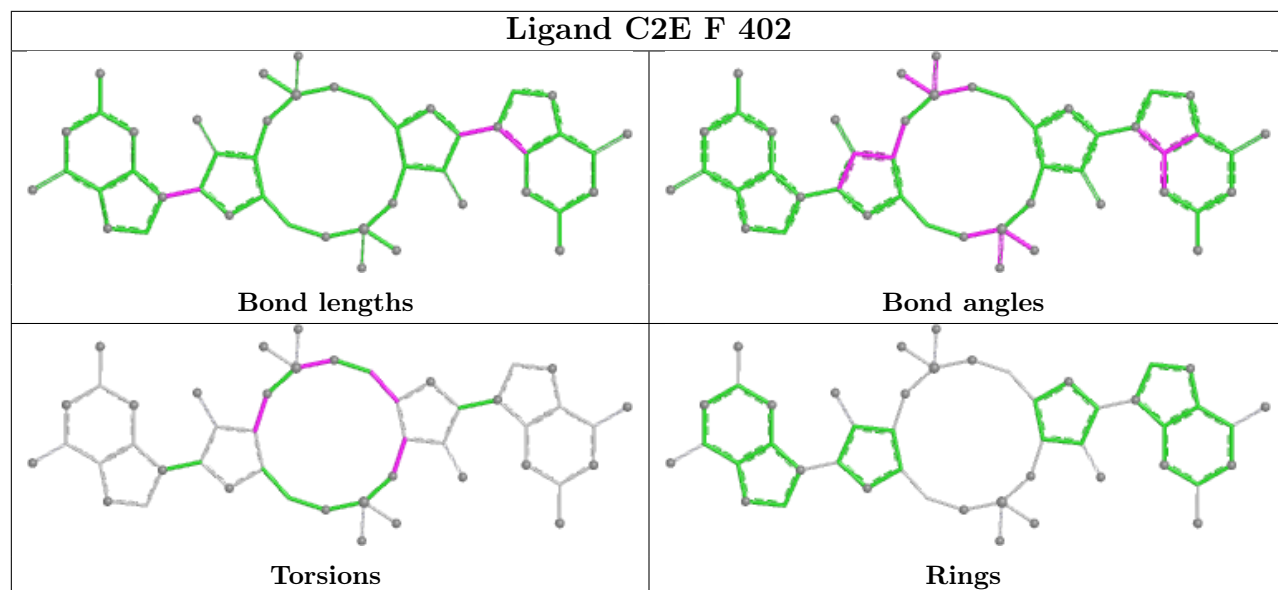


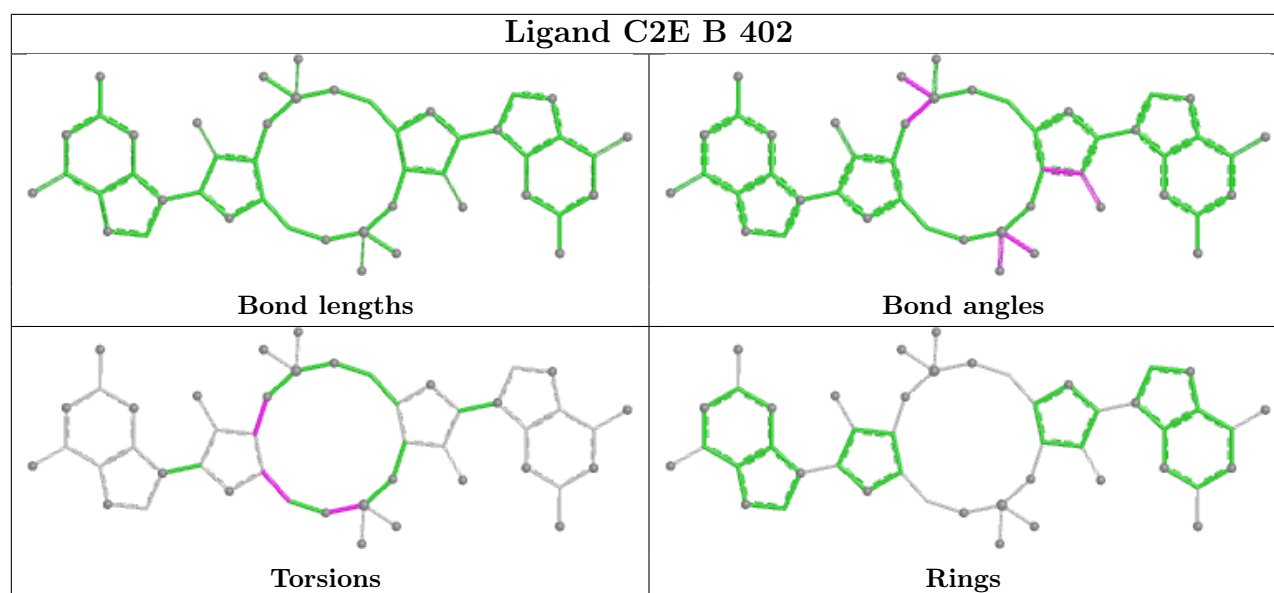
Ligand C2E C 402



Ligand C2E D 402







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	286:ALA	C	287:GLN	N	2.50

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	298/308 (96%)	-0.25	1 (0%) 90 86	43, 73, 113, 135	0
1	B	298/308 (96%)	-0.03	6 (2%) 65 56	52, 84, 147, 184	0
1	C	281/308 (91%)	-0.15	11 (3%) 43 34	41, 68, 128, 161	0
1	D	281/308 (91%)	-0.07	10 (3%) 46 37	48, 81, 137, 165	0
1	E	281/308 (91%)	-0.20	4 (1%) 73 64	45, 75, 129, 170	0
1	F	281/308 (91%)	0.23	8 (2%) 55 45	54, 98, 134, 157	0
All	All	1720/1848 (93%)	-0.08	40 (2%) 61 51	41, 80, 132, 184	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	284	PHE	5.4
1	C	287	GLN	4.0
1	B	292	ALA	3.9
1	D	284	PHE	3.7
1	E	198	VAL	3.7
1	A	198	VAL	3.4
1	C	169	TRP	3.4
1	D	198	VAL	3.2
1	B	198	VAL	3.2
1	F	190	PHE	3.1
1	D	169	TRP	3.1
1	E	206	PRO	3.0
1	C	205	LYS	2.8
1	F	198	VAL	2.7
1	D	216	LEU	2.7
1	C	290	GLU	2.6
1	D	287	GLN	2.6
1	F	276	CYS	2.6
1	D	277	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	198	VAL	2.5
1	C	220	GLY	2.4
1	F	271	LEU	2.4
1	F	195	LEU	2.4
1	D	222	TYR	2.4
1	B	305	HIS	2.4
1	B	216	LEU	2.3
1	F	169	TRP	2.3
1	E	284	PHE	2.3
1	F	24	PRO	2.3
1	F	205	LYS	2.3
1	B	283	ARG	2.2
1	D	219	VAL	2.2
1	E	273	THR	2.2
1	C	216	LEU	2.1
1	B	273	THR	2.1
1	D	279	ILE	2.1
1	C	275	VAL	2.0
1	C	206	PRO	2.0
1	D	283	ARG	2.0
1	C	273	THR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MG	F	403	1/1	0.92	0.13	55,55,55,55	0
2	C2E	D	401	46/46	0.93	0.08	47,53,69,82	0

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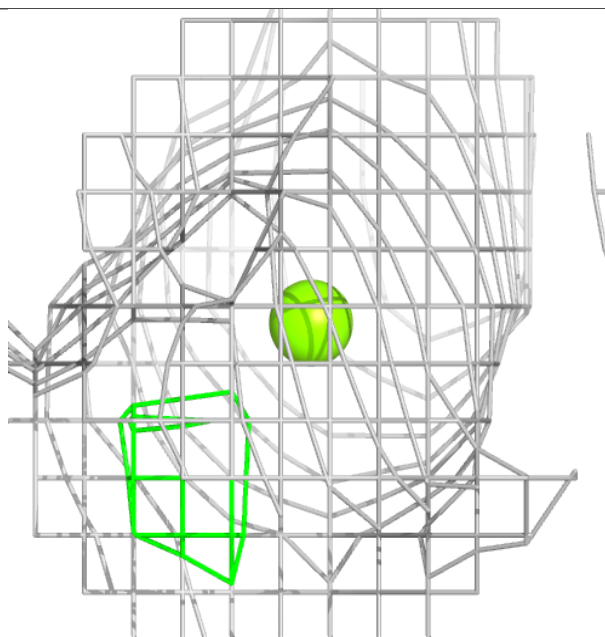
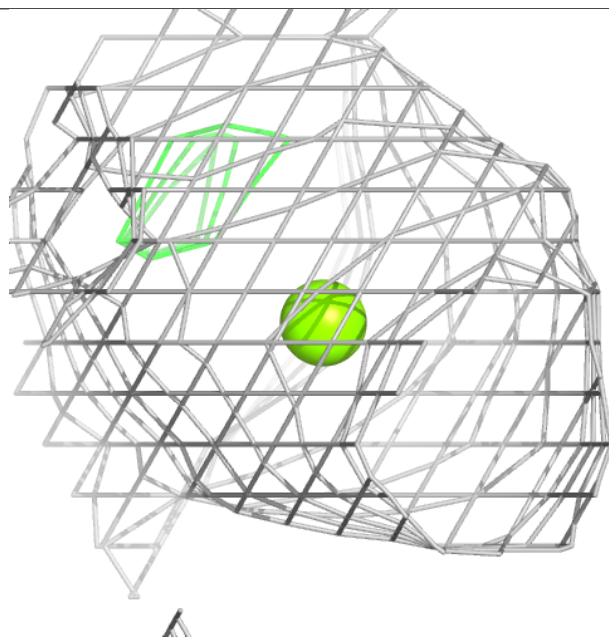
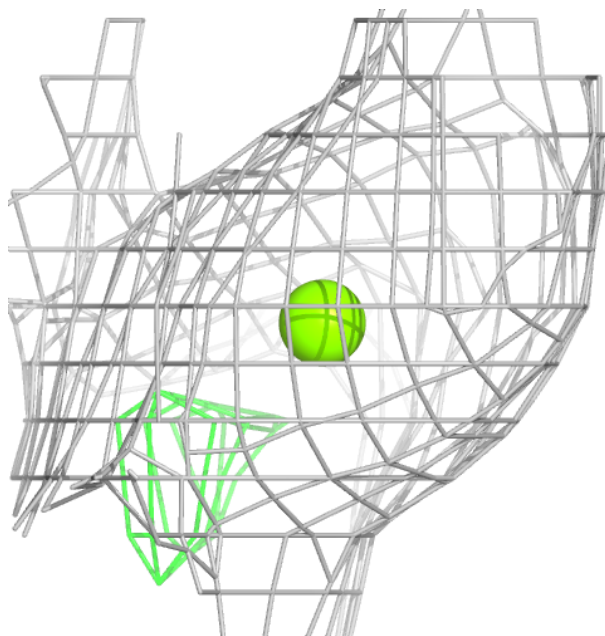
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	C2E	D	402	46/46	0.93	0.09	52,60,66,75	0
2	C2E	C	402	46/46	0.93	0.08	49,56,59,62	0
2	C2E	C	401	46/46	0.94	0.08	50,56,72,75	0
3	MG	B	403	1/1	0.95	0.10	53,53,53,53	0
3	MG	A	403	1/1	0.95	0.09	39,39,39,39	0
3	MG	C	403	1/1	0.96	0.08	39,39,39,39	0
2	C2E	E	401	46/46	0.96	0.06	49,57,71,73	0
2	C2E	B	401	46/46	0.97	0.05	48,55,59,63	0
2	C2E	F	401	46/46	-	-	33,45,58,59	46
2	C2E	F	402	46/46	-	-	43,53,58,64	46
3	MG	E	403	1/1	0.97	0.09	44,44,44,44	0
2	C2E	E	402	46/46	0.97	0.05	52,62,67,70	0
2	C2E	B	402	46/46	0.98	0.04	49,54,57,59	0
2	C2E	A	402	46/46	0.98	0.04	45,52,56,57	0
2	C2E	A	401	46/46	0.98	0.04	44,50,56,58	0
3	MG	D	403	1/1	0.99	0.04	59,59,59,59	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.

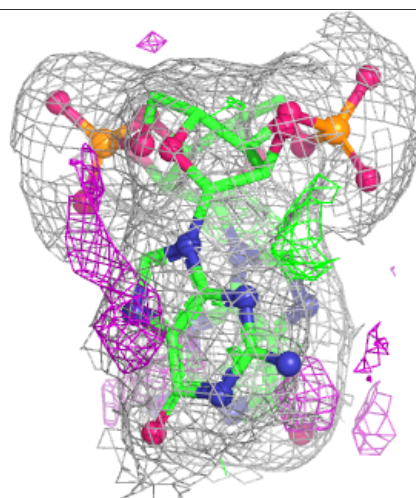
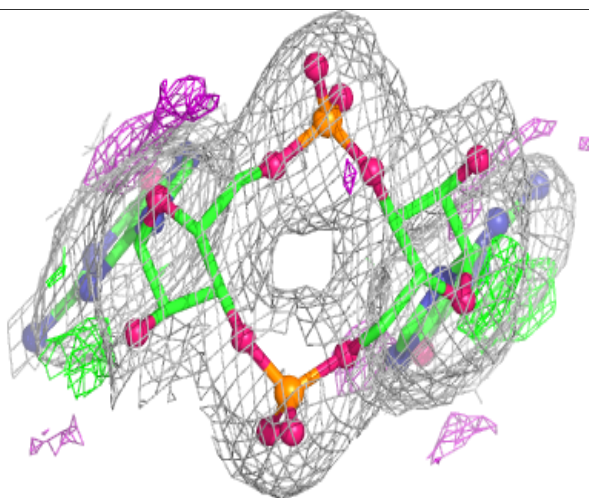
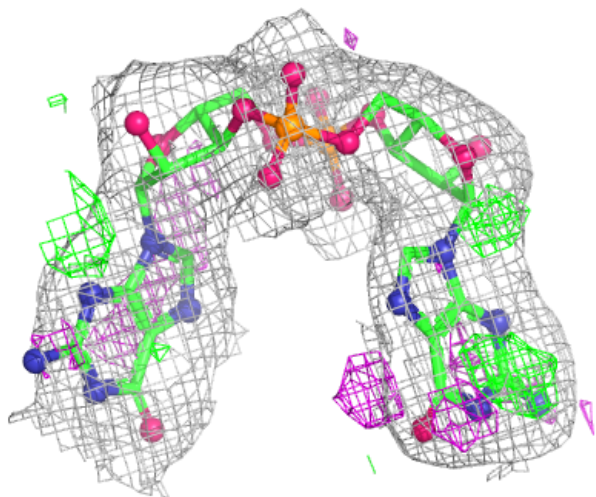
Electron density around MG F 403:

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



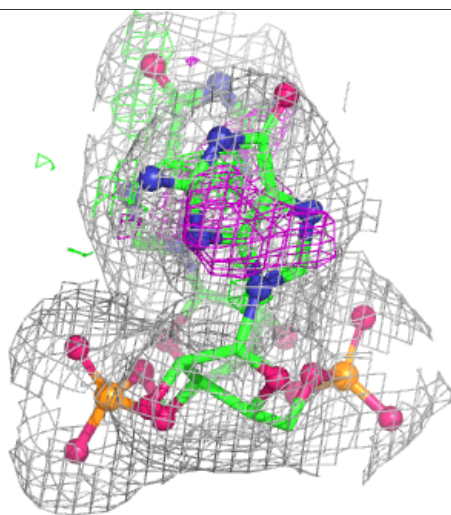
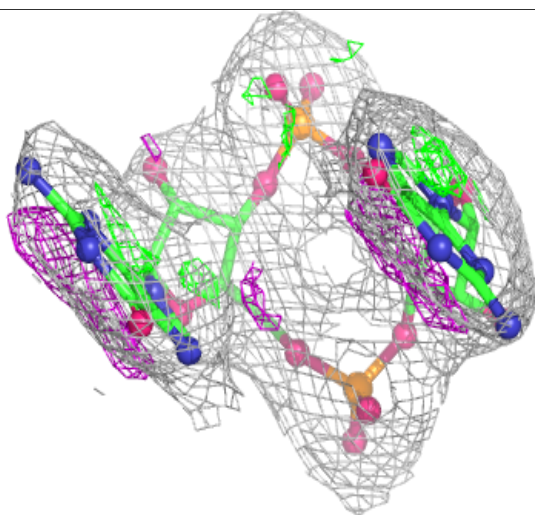
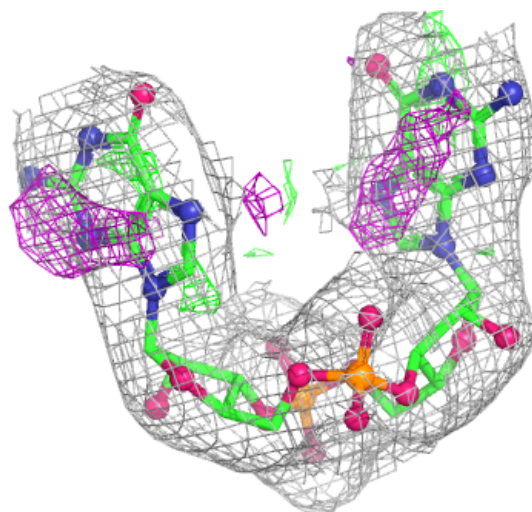
Electron density around C2E D 401:

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and green (positive)



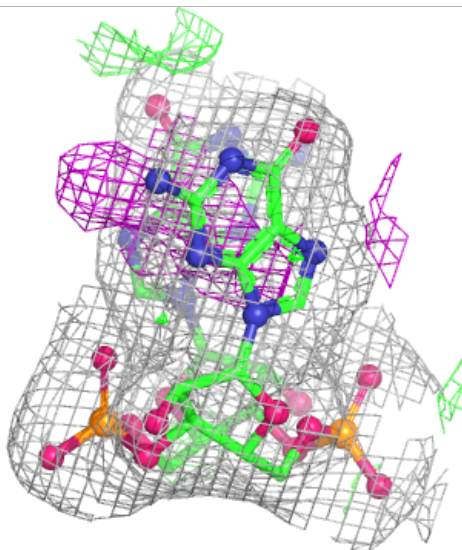
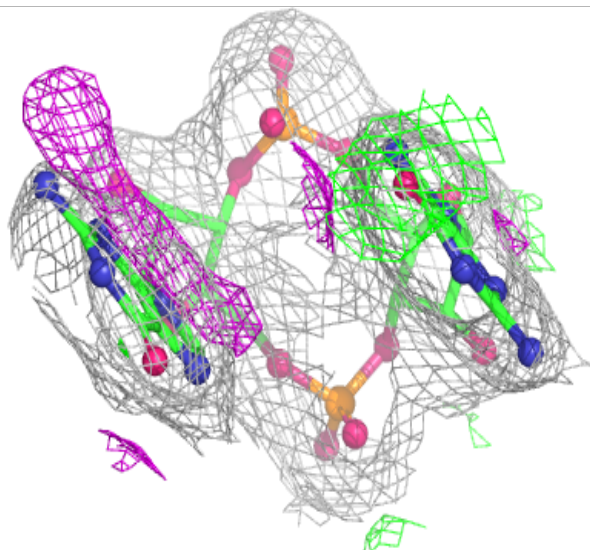
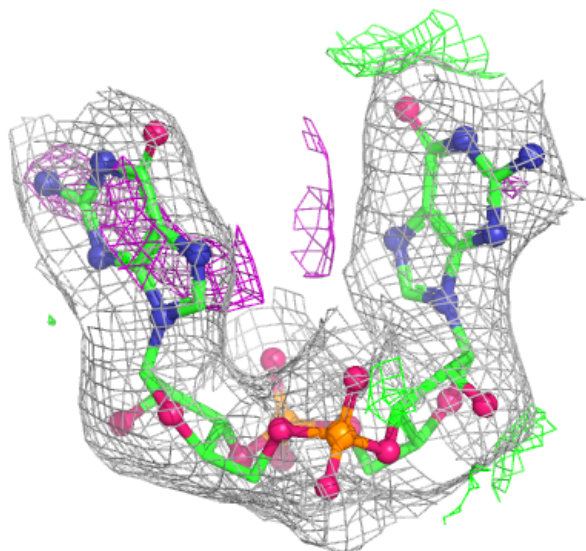
Electron density around C2E D 402:

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



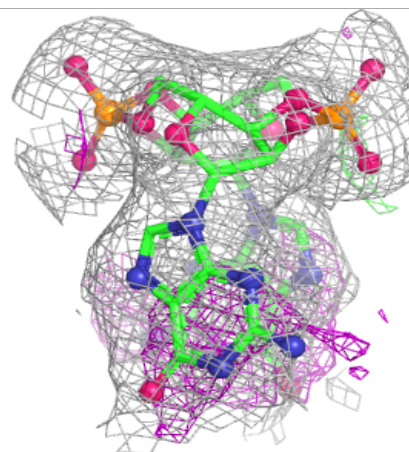
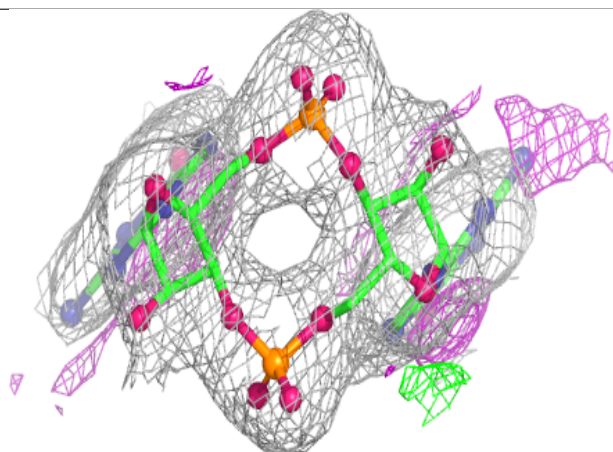
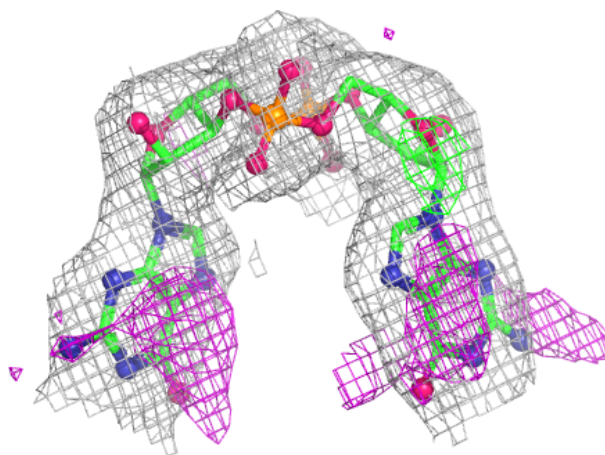
Electron density around C2E C 402:

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and green (positive)



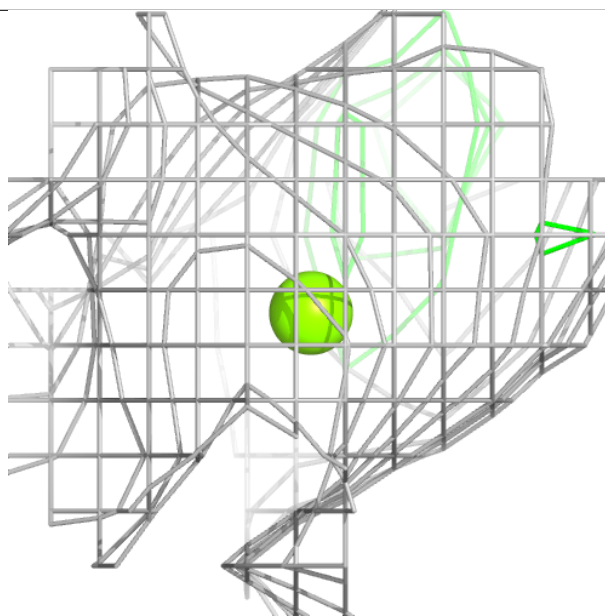
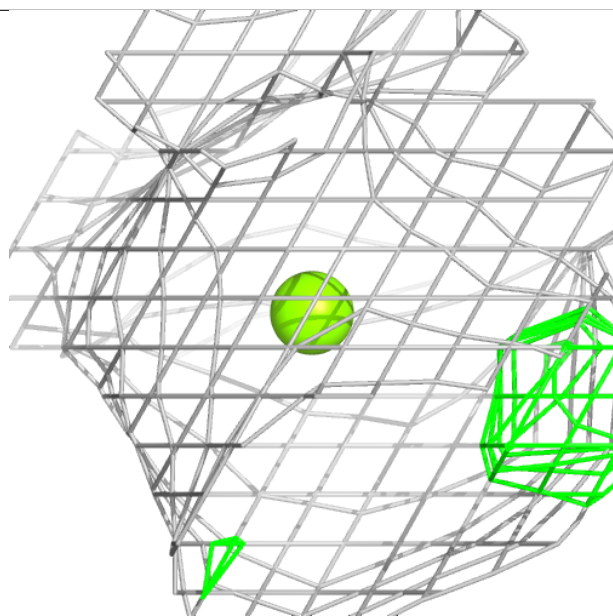
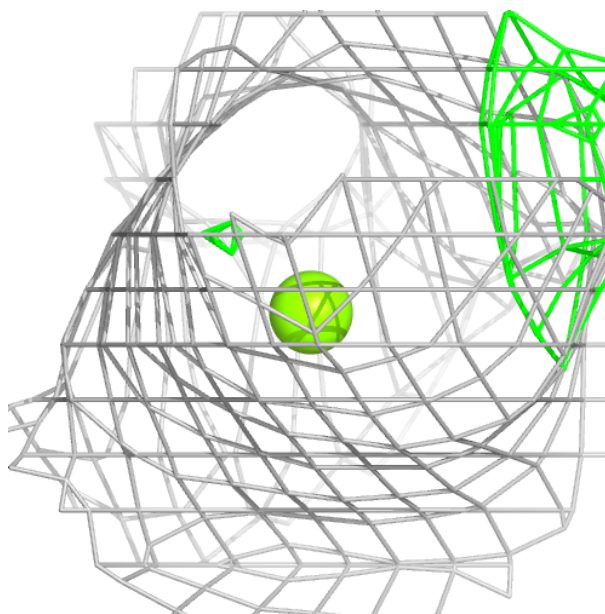
Electron density around C2E C 401:

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and green (positive)



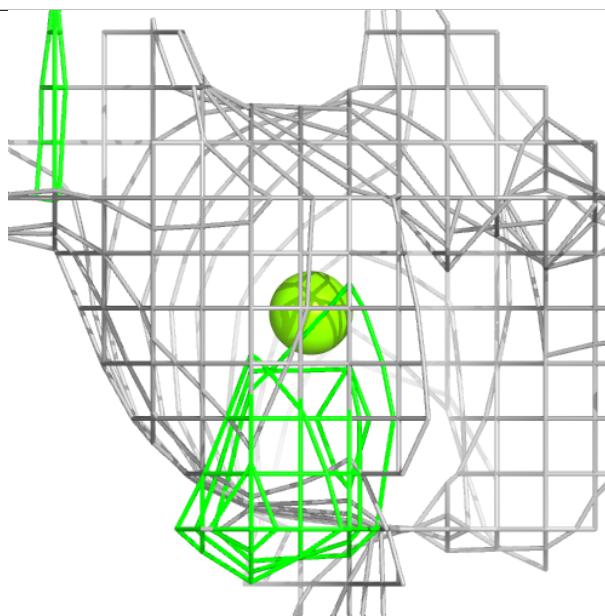
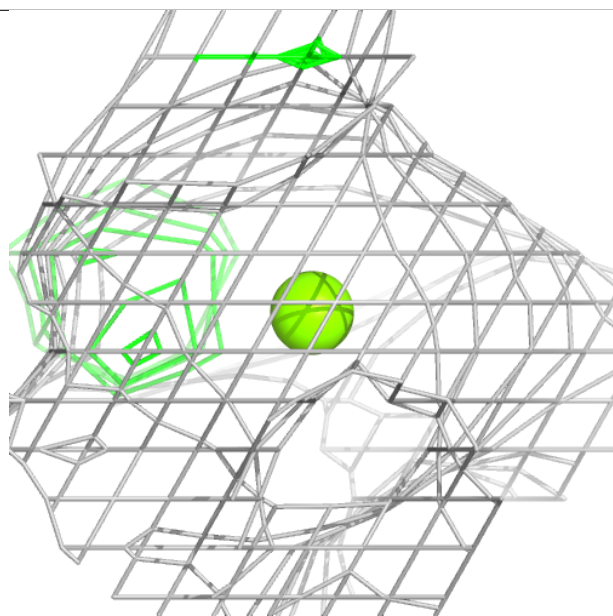
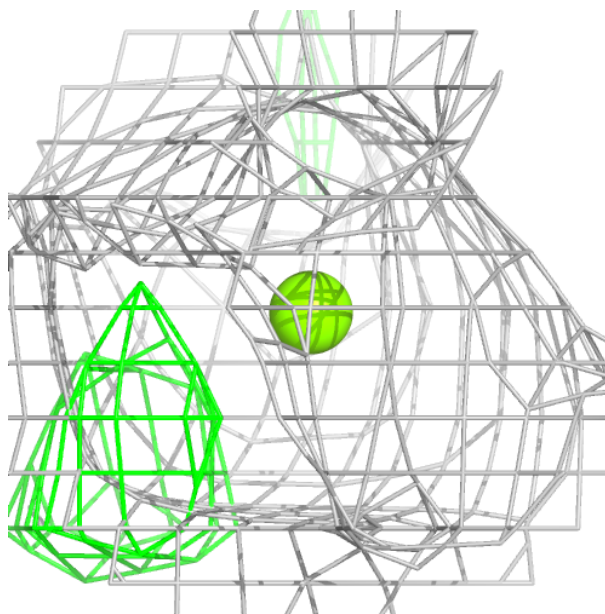
Electron density around MG B 403:

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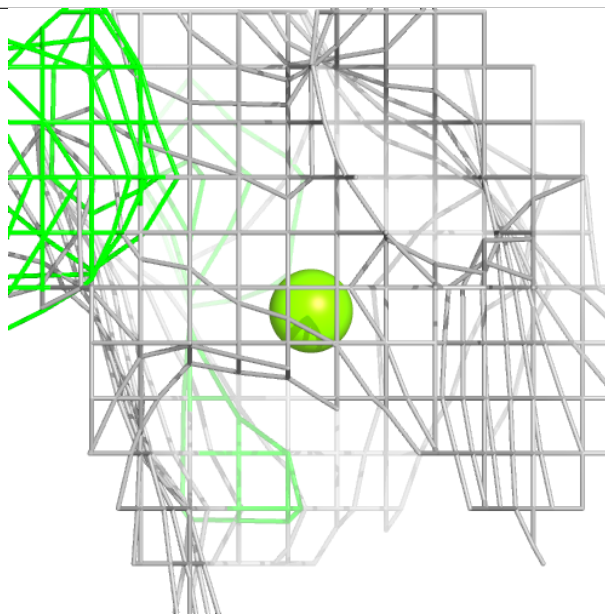
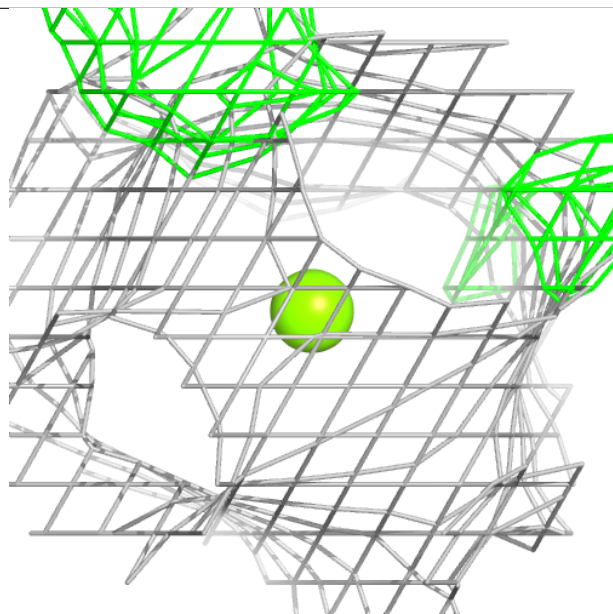
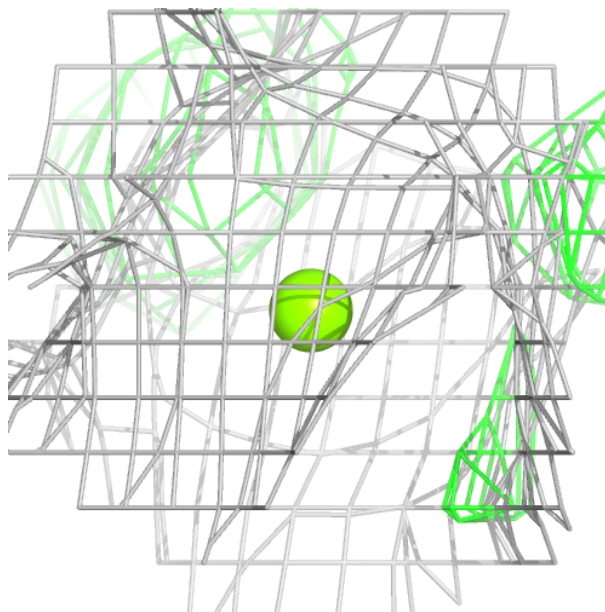
Electron density around MG A 403:

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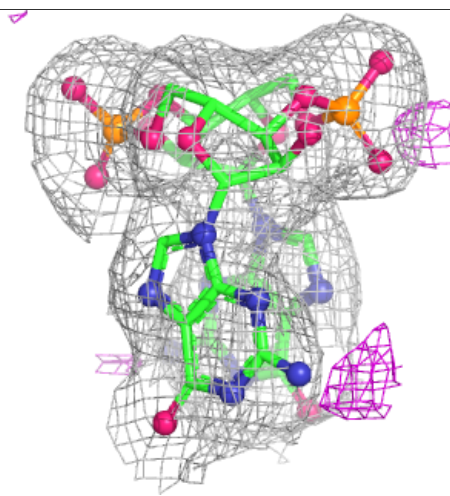
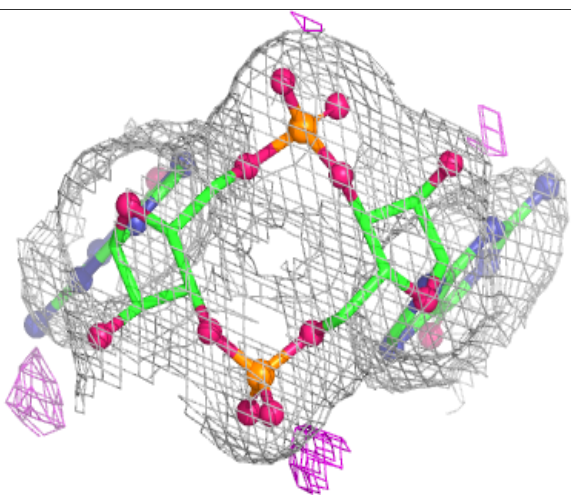
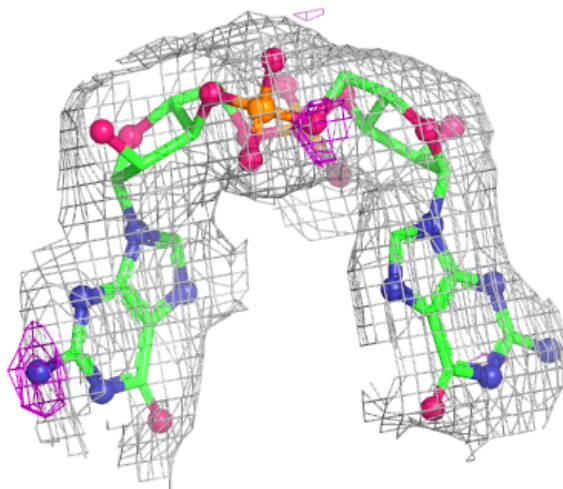
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and green (positive)



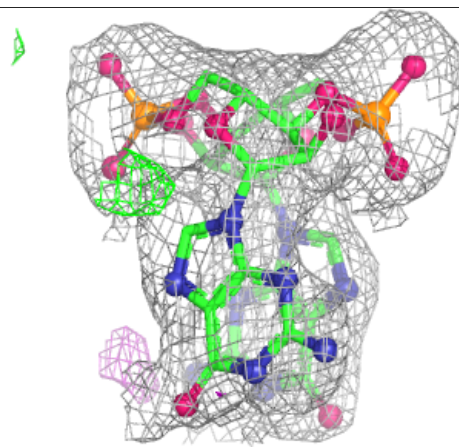
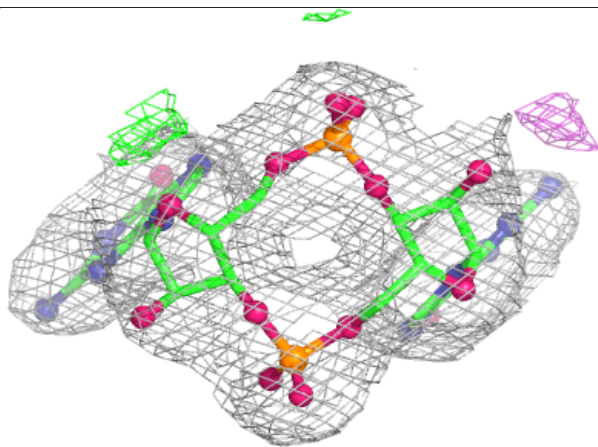
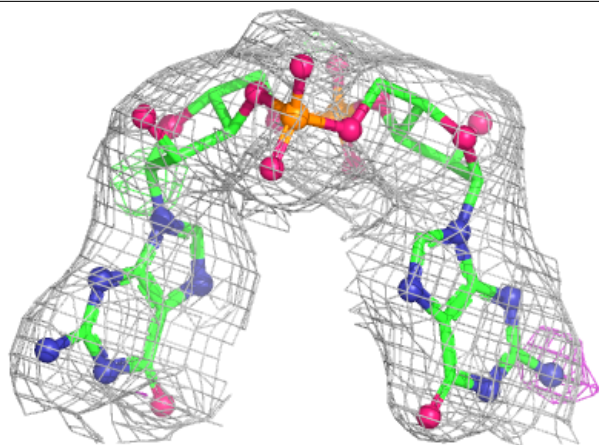
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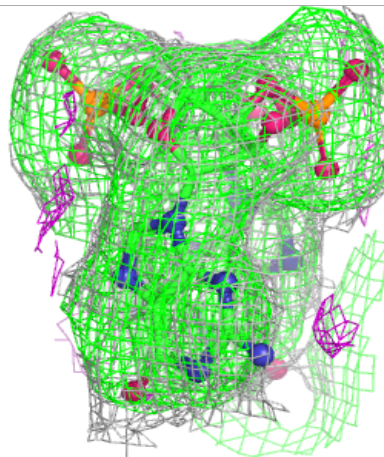
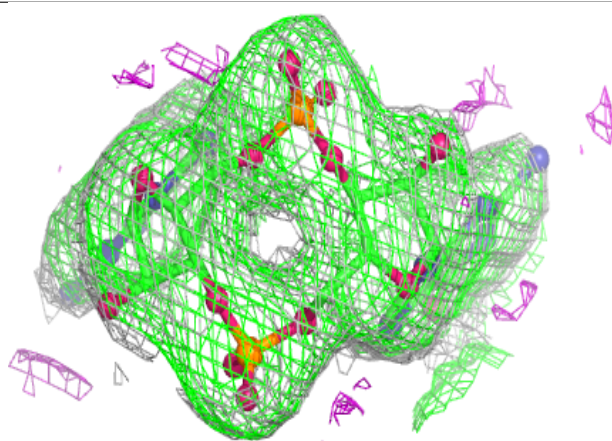
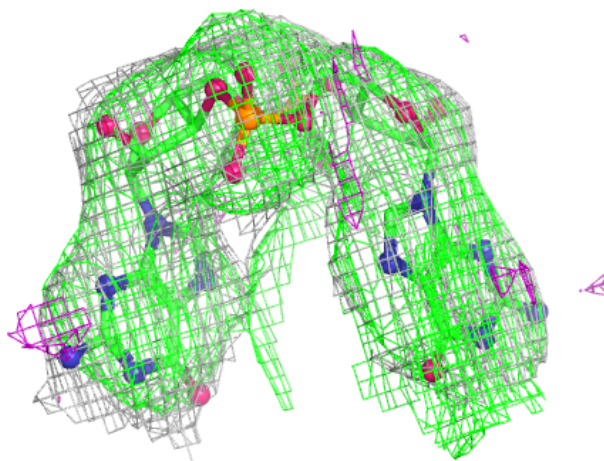
Electron density around C2E B 401:

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and green (positive)



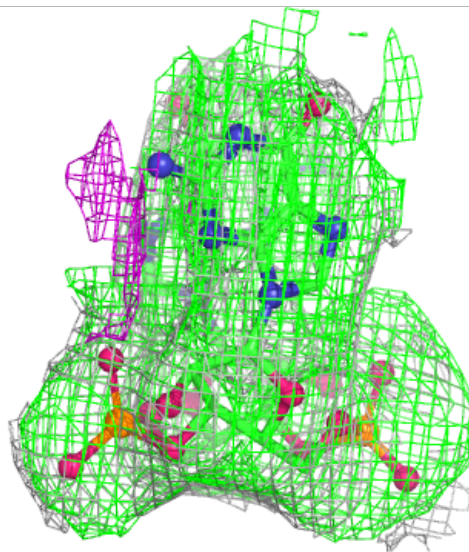
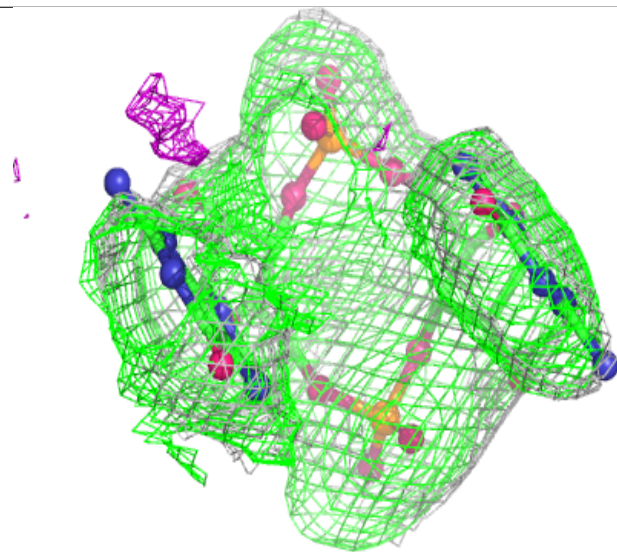
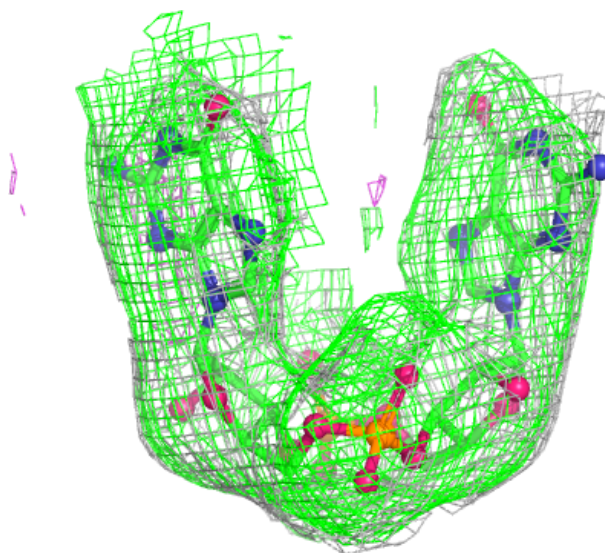
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and green (positive)



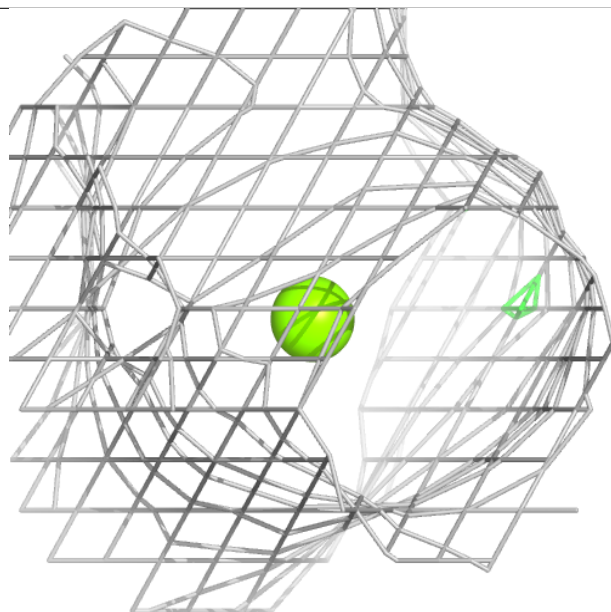
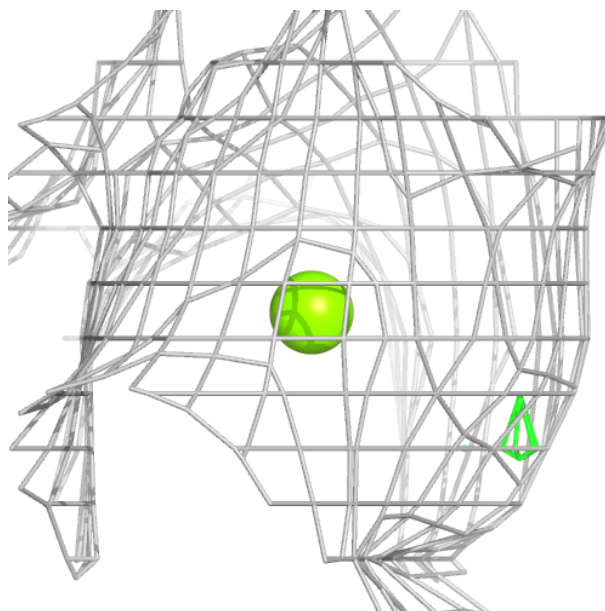
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and green (positive)



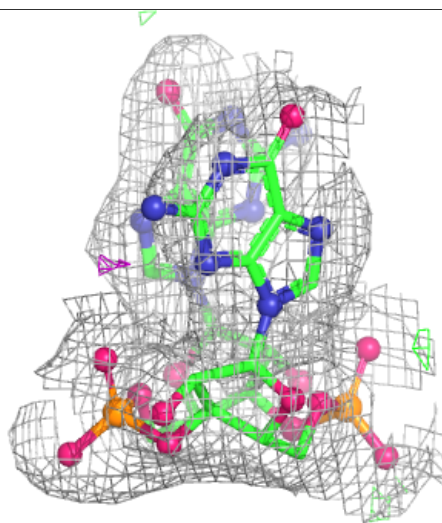
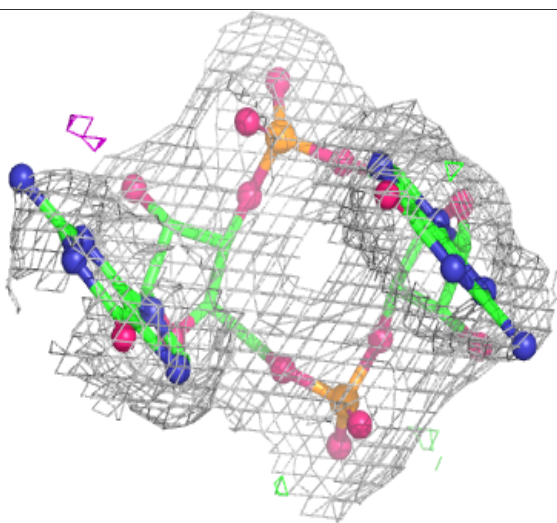
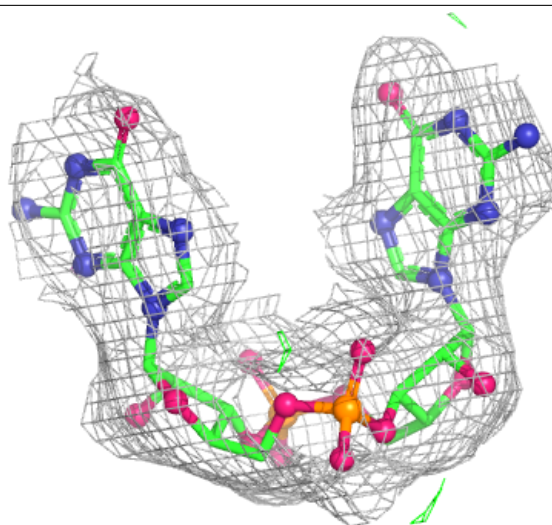
Electron density around MG E 403:

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and green (positive)



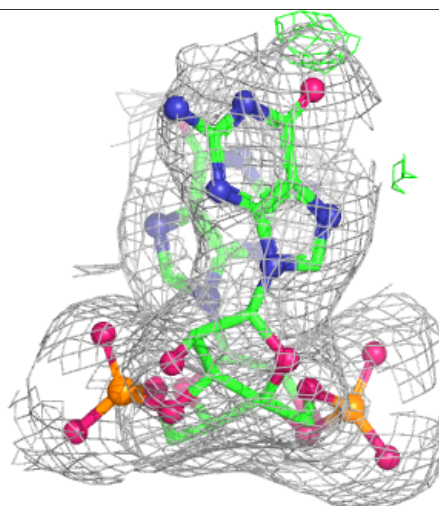
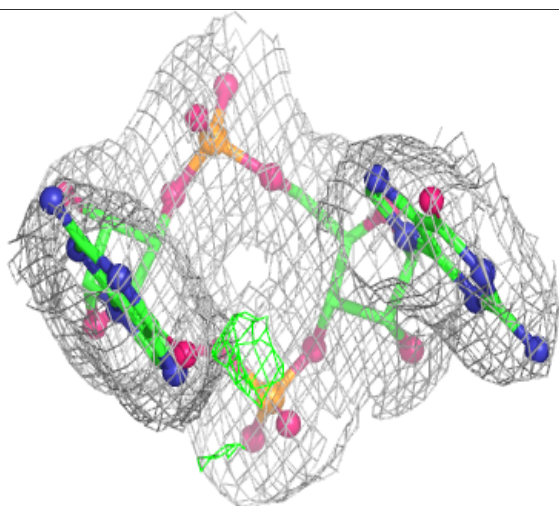
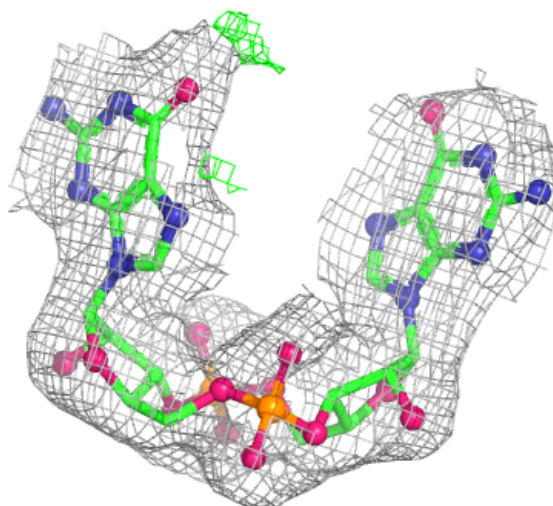
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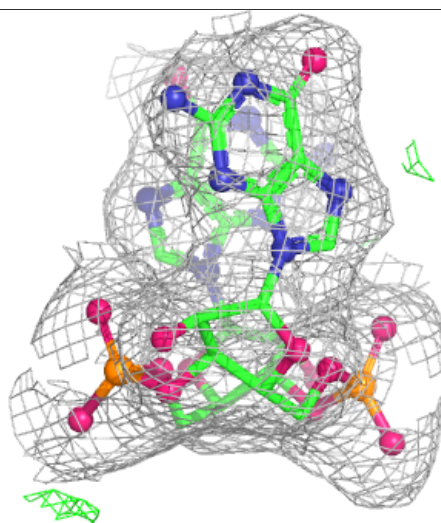
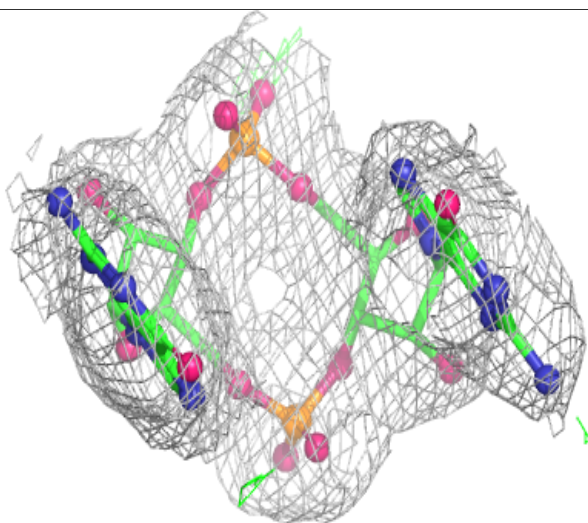
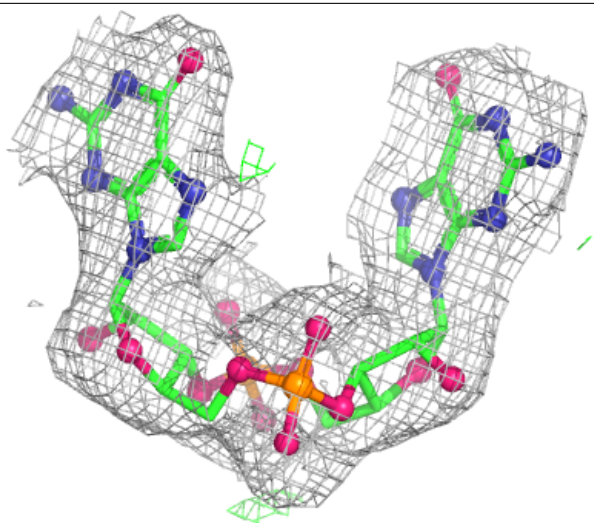
Electron density around C2E B 402:

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and green (positive)



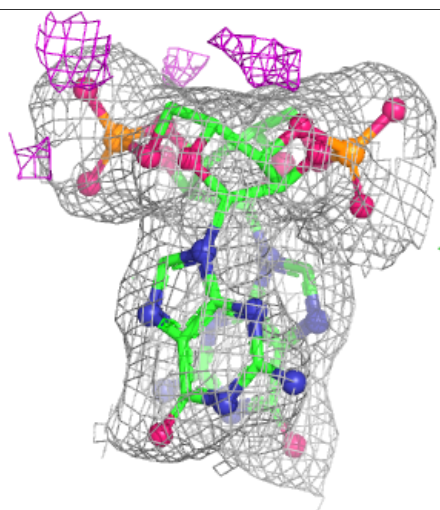
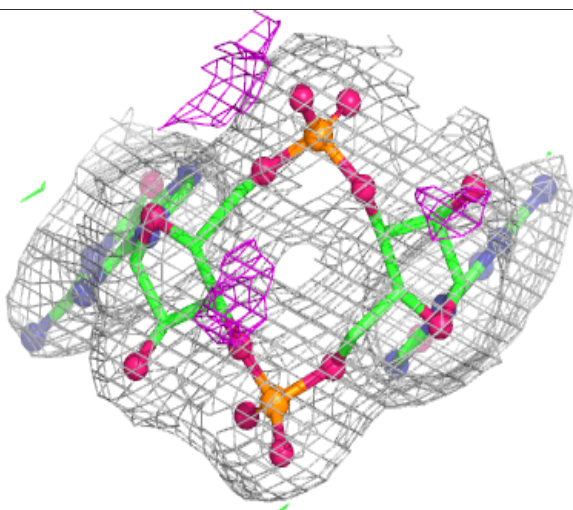
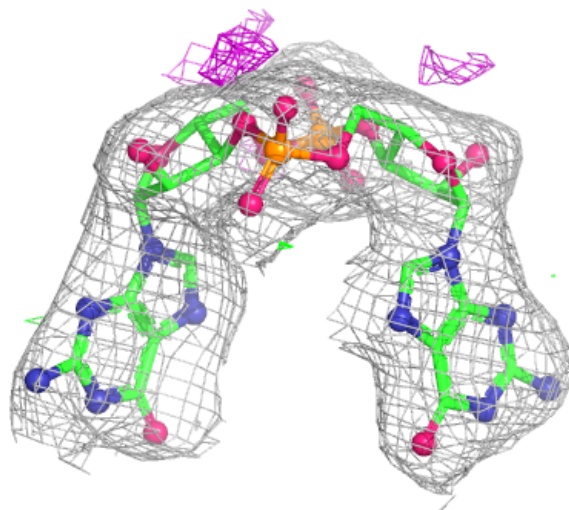
Electron density around C2E A 402:

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and green (positive)



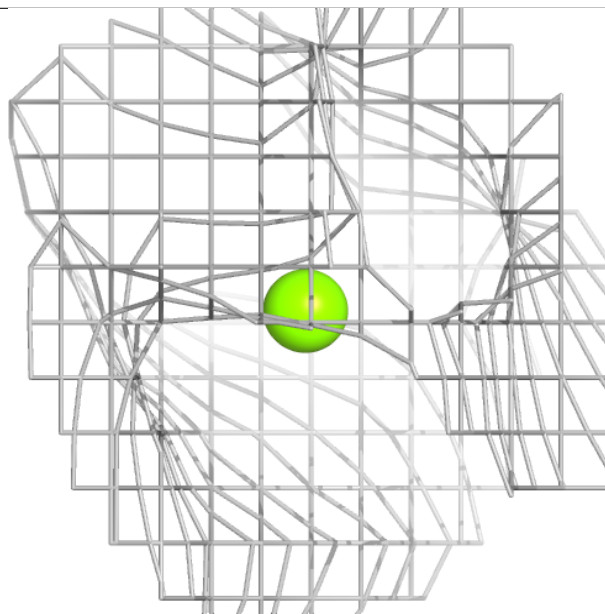
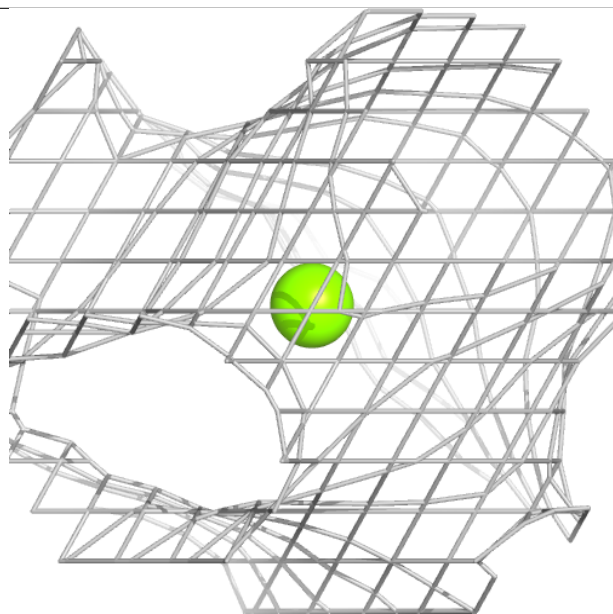
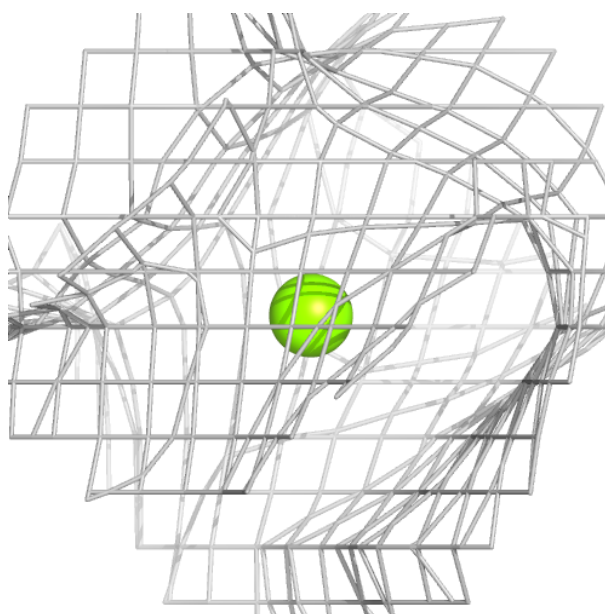
Electron density around C2E A 401:

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



Electron density around MG D 403:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.