



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:53 PM UTC

PDB ID : 4KL1 / pdb_00004kl1
Title : HCN4 CNBD in complex with cGMP
Authors : Lolicato, M.; Arrigoni, C.; Zucca, S.; Nardini, M.; Bucchi, A.; Schroeder, I.; Simmons, K.; Bolognesi, M.; DiFrancesco, D.; Schwede, F.; Fishwick, C.W.G.; Johnson, A.P.K.; Thiel, G.; Moroni, A.
Deposited on : 2013-05-07
Resolution : 2.70 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

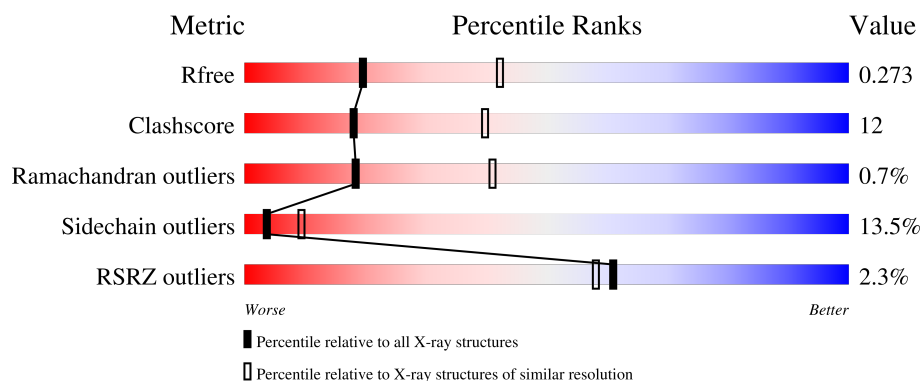
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	199	2% 67% 22% 7% . .
1	B	199	4% 61% 28% 7% . .
1	C	199	% 68% 25% . .
1	D	199	3% 60% 29% 7% . .

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ACT	A	803	-	-	X	-
3	ACT	A	804	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6814 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 Potassium/sodium hyperpolarization-activated cyclic nucleotide-gated channel 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	0	2	0
			1615	1021	284	299	11			
1	B	193	Total	C	N	O	S	0	6	0
			1637	1034	287	306	10			
1	C	193	Total	C	N	O	S	0	3	0
			1624	1026	287	301	10			
1	D	193	Total	C	N	O	S	0	1	0
			1607	1015	281	301	10			

There are 24 discrepancies between the modelled and reference sequences:

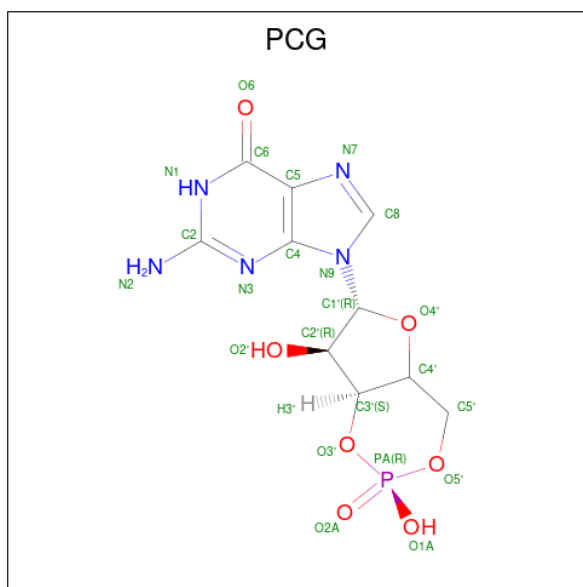
Chain	Residue	Modelled	Actual	Comment	Reference
A	515	GLY	-	expression tag	UNP Q9Y3Q4
A	516	PRO	-	expression tag	UNP Q9Y3Q4
A	517	SER	-	expression tag	UNP Q9Y3Q4
A	518	SER	-	expression tag	UNP Q9Y3Q4
A	519	PRO	-	expression tag	UNP Q9Y3Q4
A	520	MET	-	expression tag	UNP Q9Y3Q4
B	515	GLY	-	expression tag	UNP Q9Y3Q4
B	516	PRO	-	expression tag	UNP Q9Y3Q4
B	517	SER	-	expression tag	UNP Q9Y3Q4
B	518	SER	-	expression tag	UNP Q9Y3Q4
B	519	PRO	-	expression tag	UNP Q9Y3Q4
B	520	MET	-	expression tag	UNP Q9Y3Q4
C	515	GLY	-	expression tag	UNP Q9Y3Q4
C	516	PRO	-	expression tag	UNP Q9Y3Q4
C	517	SER	-	expression tag	UNP Q9Y3Q4
C	518	SER	-	expression tag	UNP Q9Y3Q4
C	519	PRO	-	expression tag	UNP Q9Y3Q4
C	520	MET	-	expression tag	UNP Q9Y3Q4
D	515	GLY	-	expression tag	UNP Q9Y3Q4
D	516	PRO	-	expression tag	UNP Q9Y3Q4

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Chain	Residue	Modelled	Actual	Comment	Reference
D	517	SER	-	expression tag	UNP Q9Y3Q4
D	518	SER	-	expression tag	UNP Q9Y3Q4
D	519	PRO	-	expression tag	UNP Q9Y3Q4
D	520	MET	-	expression tag	UNP Q9Y3Q4

- Molecule 2 is CYCLIC GUANOSINE MONOPHOSPHATE (CCD ID: PCG) (formula: $C_{10}H_{12}N_5O_7P$).



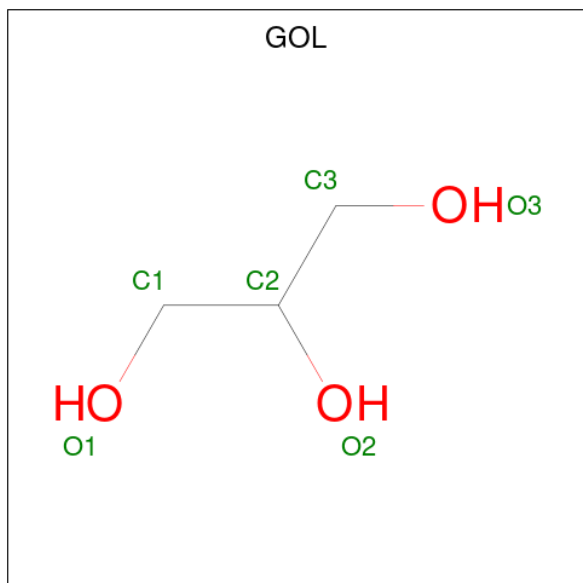
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	C	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	C	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	D	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	D	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		

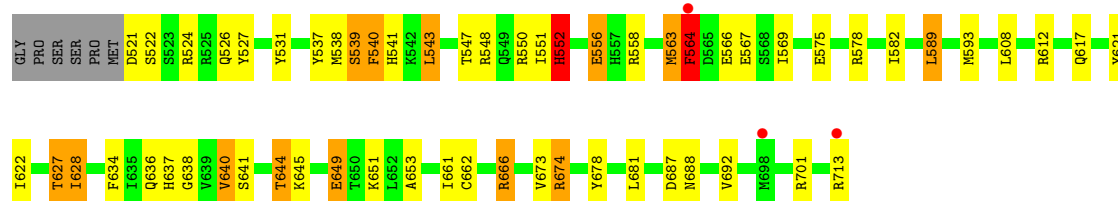
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	40	Total	O	0	0
			40	40		
5	B	17	Total	O	0	0
			17	17		
5	C	43	Total	O	0	0
			43	43		
5	D	18	Total	O	0	1
			19	19		

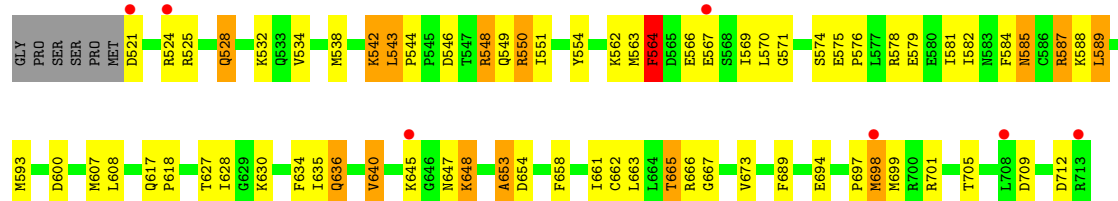
3 Residue-property plots [i](#)

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: Potassium/sodium hyperpolarization-activated cyclic nucleotide-gated channel 4



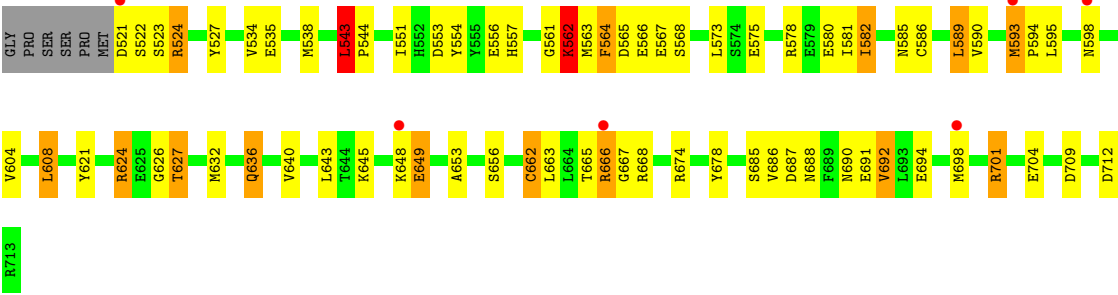
- Molecule 1: Potassium/sodium hyperpolarization-activated cyclic nucleotide-gated channel 4



- Molecule 1: Potassium/sodium hyperpolarization-activated cyclic nucleotide-gated channel 4



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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	97.10Å 99.03Å 109.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	58.62 – 2.70 58.62 – 2.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (58.62-2.70) 100.0 (58.62-2.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.73 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.210 , 0.272 0.206 , 0.273	Depositor DCC
R_{free} test set	1505 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	41.6	Xtriage
Anisotropy	0.080	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.019 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6814	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% 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: GOL, CSO, ACT, PCG

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	1.09	2/1646 (0.1%)	1.14	8/2209 (0.4%)
1	B	1.03	2/1681 (0.1%)	1.14	5/2256 (0.2%)
1	C	1.01	1/1658 (0.1%)	1.14	5/2225 (0.2%)
1	D	1.01	3/1635 (0.2%)	1.23	12/2196 (0.5%)
All	All	1.04	8/6620 (0.1%)	1.16	30/8886 (0.3%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	636	GLN	CD-OE1	6.29	1.35	1.23
1	C	598	ASN	CG-OD1	5.71	1.34	1.23
1	A	636	GLN	CD-OE1	5.66	1.34	1.23
1	D	636	GLN	CD-NE2	-5.57	1.21	1.33
1	D	636	GLN	CD-OE1	5.32	1.33	1.23
1	A	552	HIS	ND1-CE1	5.16	1.37	1.32
1	B	635	ILE	CA-CB	-5.06	1.47	1.53
1	D	557	HIS	ND1-CE1	5.06	1.37	1.32

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	542	LYS	N-CA-C	8.60	121.77	111.02
1	D	524	ARG	N-CA-C	-8.12	102.37	111.14
1	C	522	SER	N-CA-C	-7.25	104.47	113.38
1	C	569	ILE	N-CA-C	-7.08	103.95	110.82
1	A	552	HIS	CB-CG-CD2	-6.73	122.46	131.20
1	D	543	LEU	CA-C-N	6.64	124.51	119.66
1	D	543	LEU	C-N-CA	6.64	124.51	119.66
1	D	557	HIS	CB-CG-CD2	-6.57	122.66	131.20
1	D	662	CYS	N-CA-C	6.48	118.42	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	645	LYS	N-CA-C	6.24	118.96	108.23
1	A	653	ALA	N-CA-C	5.99	115.92	108.19
1	D	522	SER	N-CA-C	-5.98	106.32	113.97
1	B	653	ALA	N-CA-C	5.92	115.92	108.45
1	D	653	ALA	N-CA-C	5.90	116.21	108.07
1	A	564	PHE	N-CA-C	5.82	118.54	109.52
1	D	594	PRO	N-CA-C	5.81	124.44	112.47
1	A	552	HIS	CB-CG-ND1	5.72	131.28	122.70
1	B	635	ILE	N-CA-C	5.65	115.69	108.12
1	D	557	HIS	CB-CG-ND1	5.61	131.11	122.70
1	A	522	SER	N-CA-C	-5.44	105.78	112.90
1	D	562	LYS	N-CA-C	5.35	117.55	108.02
1	D	580	GLU	N-CA-C	-5.33	105.55	111.36
1	A	678	TYR	CA-C-N	-5.14	115.75	122.99
1	A	678	TYR	C-N-CA	-5.14	115.75	122.99
1	A	687	ASP	N-CA-C	-5.10	105.61	111.07
1	D	692	VAL	N-CA-C	5.06	115.83	110.72
1	B	564	PHE	N-CA-C	5.03	115.95	108.86
1	C	600	ASP	CA-C-N	5.01	126.11	119.84
1	C	600	ASP	C-N-CA	5.01	126.11	119.84
1	B	549	GLN	N-CA-C	-5.00	105.53	111.69

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1615	0	1596	43	0
1	B	1637	0	1617	52	0
1	C	1624	0	1606	38	0
1	D	1607	0	1578	43	0
2	A	46	0	22	1	0
2	B	46	0	22	2	0
2	C	46	0	22	1	0
2	D	46	0	22	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	12	0	9	5	0
3	C	4	0	3	0	0
4	C	12	0	16	3	0
5	A	40	0	0	0	0
5	B	17	0	0	0	0
5	C	43	0	0	3	0
5	D	19	0	0	0	0
All	All	6814	0	6513	161	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 12.

All (161) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:593:MET:HE1	1:A:634:PHE:CD2	1.98	0.99
1:A:593:MET:HE1	1:A:634:PHE:HD2	1.29	0.95
1:A:674:ARG:HH11	1:A:674:ARG:HG2	1.38	0.88
1:D:538:MET:HA	1:D:543:LEU:HD22	1.64	0.80
1:B:593:MET:HE1	1:B:634:PHE:CD2	2.17	0.79
1:D:586:CSO:O	1:D:590:VAL:HG23	1.85	0.76
3:A:803:ACT:H2	1:C:624[A]:ARG:NH1	2.01	0.75
1:A:688:ASN:O	1:A:692:VAL:HG23	1.86	0.75
1:B:587:ARG:HH11	1:B:587:ARG:CG	2.01	0.74
1:D:604:VAL:O	1:D:608:LEU:HD22	1.89	0.73
1:D:582:ILE:CG2	1:D:608:LEU:HB3	2.19	0.73
1:A:638:GLY:HA2	3:A:804:ACT:H2	1.70	0.72
1:B:587:ARG:HH11	1:B:587:ARG:HG2	1.55	0.71
1:B:593:MET:HE1	1:B:634:PHE:HD2	1.56	0.69
1:C:567:GLU:OE1	1:C:578[B]:ARG:NH2	2.26	0.69
1:C:713:ARG:O	1:C:713:ARG:HG3	1.91	0.68
1:B:627:THR:HG22	1:B:628:ILE:N	2.09	0.68
1:A:531:TYR:OH	1:A:552:HIS:HB2	1.94	0.68
1:A:527:TYR:HA	1:A:563:MET:HG3	1.79	0.64
1:B:524:ARG:O	1:B:528:GLN:HG2	1.98	0.64
1:A:593:MET:CE	1:A:634:PHE:HD2	2.06	0.63
1:B:640:VAL:CG1	1:B:673:VAL:HG13	2.29	0.63
1:D:521:ASP:HB3	1:D:562:LYS:NZ	2.14	0.63
1:B:538:MET:HG2	1:B:543:LEU:CD2	2.27	0.63
4:C:804:GOL:H32	1:D:624:ARG:HH12	1.64	0.62
1:A:575:GLU:OE1	1:A:578[A]:ARG:NH1	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:632:MET:HB3	1:C:663:LEU:HD22	1.81	0.62
1:B:648:LYS:N	1:B:648:LYS:HD3	2.14	0.62
1:D:521:ASP:HB3	1:D:562:LYS:HZ3	1.65	0.62
1:B:593:MET:CE	1:B:634:PHE:HD2	2.13	0.61
1:A:649:GLU:OE1	1:A:651:LYS:NZ	2.34	0.60
1:A:640:VAL:HG13	1:A:673:VAL:HG13	1.83	0.60
1:A:637:HIS:NE2	3:A:804:ACT:H3	2.17	0.60
1:B:538:MET:HG2	1:B:543:LEU:HD23	1.83	0.60
1:D:632:MET:HB3	1:D:663:LEU:HD22	1.83	0.59
1:A:537:TYR:HE1	1:B:566[B]:GLU:HG3	1.68	0.59
1:B:538:MET:SD	1:B:548:ARG:HG2	2.42	0.59
1:D:624:ARG:O	1:D:627:THR:OG1	2.21	0.58
4:C:804:GOL:H31	1:D:621:TYR:O	2.03	0.58
1:A:564:PHE:HD2	1:A:564:PHE:H	1.51	0.58
1:A:538:MET:HA	1:A:543:LEU:HD22	1.86	0.58
1:C:593:MET:O	1:C:594:PRO:C	2.42	0.58
1:A:537:TYR:HE1	1:B:566[B]:GLU:CG	2.16	0.58
1:B:665:THR:O	1:B:667:GLY:N	2.37	0.57
1:D:688:ASN:O	1:D:692:VAL:HG23	2.03	0.57
1:B:694:GLU:O	1:B:697:PRO:HD3	2.05	0.57
1:C:521:ASP:HB2	1:C:524:ARG:HB2	1.86	0.56
1:A:674:ARG:HG2	1:A:674:ARG:NH1	2.14	0.56
1:C:624[B]:ARG:NH2	5:C:912:HOH:O	2.38	0.56
1:C:665:THR:O	1:C:666:ARG:C	2.48	0.56
1:A:564:PHE:N	1:A:564:PHE:CD2	2.75	0.55
1:D:575:GLU:OE1	1:D:578:ARG:NH1	2.37	0.55
1:A:564:PHE:HZ	1:C:540:PHE:CD2	2.23	0.55
1:D:626:GLY:O	1:D:668:ARG:HD3	2.06	0.55
1:A:637:HIS:CE1	3:A:804:ACT:H3	2.41	0.55
1:C:584:PHE:CZ	1:D:544:PRO:HD3	2.42	0.55
1:D:553:ASP:HB3	1:D:678:TYR:OH	2.07	0.55
1:A:537:TYR:HE1	1:B:566[A]:GLU:CD	2.15	0.54
1:A:564:PHE:HD2	1:A:564:PHE:N	2.06	0.54
1:B:593:MET:HE1	1:B:634:PHE:CE2	2.44	0.53
1:C:643:LEU:HD23	1:C:649:GLU:HB3	1.89	0.53
1:C:665:THR:O	1:C:665:THR:OG1	2.26	0.53
1:A:540:PHE:CD2	1:B:564:PHE:HZ	2.26	0.52
1:D:582:ILE:HG21	1:D:608:LEU:HB3	1.91	0.52
1:C:595:LEU:O	1:C:595:LEU:HG	2.09	0.52
1:A:566:GLU:CD	1:C:537:TYR:HE1	2.17	0.52
1:B:658:PHE:CE2	1:B:673:VAL:HG11	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:575:GLU:OE2	1:B:579:GLU:OE2	2.28	0.52
1:B:627:THR:CG2	1:B:628:ILE:N	2.72	0.51
1:A:713:ARG:HB2	1:A:713:ARG:NH1	2.26	0.51
1:D:685:SER:OG	1:D:688:ASN:HB2	2.10	0.51
1:D:589:LEU:O	1:D:593:MET:HG3	2.11	0.51
1:A:552:HIS:CE1	1:A:556:GLU:OE2	2.64	0.51
1:B:588:LYS:O	1:B:589:LEU:C	2.54	0.51
1:D:701:ARG:HA	1:D:704:GLU:HG3	1.93	0.51
1:D:534:VAL:O	1:D:538:MET:HG3	2.11	0.50
1:B:534:VAL:HG12	1:B:538:MET:HE2	1.94	0.50
1:B:566[A]:GLU:OE1	2:B:802:PCG:O2'	2.27	0.50
1:B:574:SER:HB2	1:B:576:PRO:HD2	1.93	0.49
1:C:521:ASP:O	1:C:525:ARG:NH1	2.45	0.49
1:A:564:PHE:HD1	5:C:921:HOH:O	1.95	0.49
1:B:617:GLN:O	1:B:618:PRO:C	2.56	0.49
1:C:553:ASP:O	1:C:557:HIS:HD2	1.95	0.49
1:C:623:ILE:HG23	1:C:669:ARG:HG3	1.94	0.49
1:D:665:THR:O	1:D:666:ARG:C	2.55	0.49
1:D:564:PHE:CD1	1:D:564:PHE:N	2.80	0.48
1:A:539:SER:C	1:A:541:HIS:H	2.21	0.48
1:C:573:LEU:HA	1:D:554:TYR:CE2	2.48	0.48
1:B:544:PRO:HB2	1:B:546[B]:ASP:OD1	2.14	0.48
1:D:521:ASP:CB	1:D:562:LYS:NZ	2.77	0.48
1:D:564:PHE:N	1:D:564:PHE:HD1	2.11	0.48
1:B:584:PHE:O	1:B:585:ASN:C	2.58	0.47
1:D:589:LEU:O	1:D:593:MET:CG	2.63	0.47
1:B:551:ILE:HD11	1:D:581:ILE:HD11	1.97	0.46
1:B:564:PHE:N	1:B:564:PHE:CD2	2.84	0.46
1:B:640:VAL:HG11	1:B:673:VAL:HG13	1.97	0.46
1:B:600:ASP:OD1	1:B:600:ASP:C	2.59	0.46
1:B:607:MET:HE3	1:B:607:MET:HB3	1.84	0.46
1:B:554:TYR:HB2	1:B:617:GLN:NE2	2.31	0.46
1:D:523:SER:HB2	1:D:563:MET:O	2.16	0.46
1:D:562:LYS:O	2:D:802:PCG:N2	2.38	0.46
1:D:686:VAL:O	1:D:687:ASP:C	2.59	0.46
1:C:647:ASN:OD1	1:C:647:ASN:N	2.49	0.45
1:A:713:ARG:HB2	1:A:713:ARG:HH11	1.82	0.45
1:C:600:ASP:C	1:C:600:ASP:OD1	2.59	0.45
1:D:565:ASP:CG	1:D:568:SER:HB2	2.41	0.45
1:A:589:LEU:O	1:A:593:MET:HG3	2.16	0.45
1:C:581:ILE:HD11	1:D:551:ILE:HD11	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:674:ARG:HH11	1:D:674:ARG:HG2	1.81	0.45
4:C:804:GOL:H32	1:D:624:ARG:NH1	2.30	0.45
1:C:566:GLU:HA	1:C:569:ILE:HD12	1.98	0.45
1:A:627:THR:CG2	1:A:628:ILE:N	2.81	0.44
1:B:587:ARG:CG	1:B:587:ARG:NH1	2.67	0.44
1:C:701:ARG:HD3	1:C:704[B]:GLU:HB3	2.00	0.44
1:A:547:THR:HA	1:A:550:ARG:NH1	2.33	0.44
1:A:551:ILE:HD11	1:B:581:ILE:HD11	1.99	0.44
1:D:690:ASN:O	1:D:691:GLU:C	2.61	0.44
1:B:570:LEU:HB3	1:B:578[A]:ARG:HG2	2.00	0.44
1:C:550:ARG:HD3	5:C:903:HOH:O	2.18	0.43
1:C:546:ASP:O	1:C:549:GLN:HB3	2.18	0.43
1:C:701:ARG:NH1	1:C:704[A]:GLU:OE1	2.52	0.43
1:B:640:VAL:HG11	1:B:673:VAL:CG1	2.49	0.43
1:C:567:GLU:CD	1:C:578[B]:ARG:HH22	2.25	0.43
1:C:671:ALA:HB3	2:C:801:PCG:H5'1	2.00	0.43
1:A:569:ILE:HG23	1:C:534:VAL:HG22	2.00	0.43
1:A:666:ARG:CG	1:A:666:ARG:HH11	2.32	0.43
1:A:641:SER:OG	1:A:651:LYS:NZ	2.44	0.43
3:A:803:ACT:H2	1:C:624[A]:ARG:CZ	2.49	0.43
1:C:612:ARG:HH11	1:C:612:ARG:HB3	1.84	0.42
1:A:550:ARG:O	1:A:617:GLN:NE2	2.39	0.42
1:A:621:TYR:CZ	1:A:674:ARG:HD3	2.54	0.42
1:B:627:THR:HG22	1:B:628:ILE:H	1.81	0.42
1:C:701:ARG:HD3	1:C:701:ARG:HA	1.55	0.42
1:D:643:LEU:HD23	1:D:649:GLU:HB2	2.02	0.42
1:B:640:VAL:CG1	1:B:673:VAL:CG1	2.98	0.42
1:A:622:ILE:HD13	1:A:681:LEU:HD11	2.02	0.41
1:A:644:THR:O	1:A:645:LYS:C	2.62	0.41
1:B:653:ALA:O	1:B:654:ASP:C	2.63	0.41
1:A:661:ILE:HD12	2:A:801:PCG:H3'	2.02	0.41
1:D:527:TYR:CD1	1:D:527:TYR:C	2.98	0.41
1:B:554:TYR:CE2	1:D:573:LEU:HA	2.56	0.41
1:B:648:LYS:HD3	1:B:648:LYS:H	1.83	0.41
1:B:698:MET:H	1:B:698:MET:HG3	1.52	0.41
1:D:665:THR:O	1:D:667:GLY:N	2.53	0.41
1:B:663:LEU:HD23	1:B:689:PHE:CG	2.55	0.41
1:C:534:VAL:O	1:C:538:MET:HG3	2.21	0.41
1:A:521:ASP:HB2	1:A:524:ARG:HB2	2.03	0.41
1:D:521:ASP:OD1	1:D:524:ARG:NH2	2.53	0.41
1:C:574:SER:HB3	1:D:554:TYR:CE1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:538:MET:HG2	1:B:543:LEU:HD22	2.01	0.41
1:C:584:PHE:CE2	1:D:544:PRO:CG	3.03	0.41
1:A:543:LEU:HD12	1:A:543:LEU:HA	1.88	0.40
1:D:561:GLY:C	1:D:562:LYS:HG2	2.46	0.40
1:B:699:MET:C	1:B:701:ARG:H	2.29	0.40
1:C:657:TYR:CD1	1:C:657:TYR:C	2.99	0.40
1:B:661:ILE:HD12	2:B:801:PCG:H3'	2.02	0.40
1:C:712:ASP:OD1	1:C:712:ASP:C	2.63	0.40
1:B:569:ILE:C	1:B:571:GLY:N	2.78	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	192/199 (96%)	186 (97%)	5 (3%)	1 (0%)	24	48
1	B	196/199 (98%)	175 (89%)	17 (9%)	4 (2%)	6	16
1	C	193/199 (97%)	188 (97%)	5 (3%)	0	100	100
1	D	191/199 (96%)	177 (93%)	13 (7%)	1 (0%)	24	48
All	All	772/796 (97%)	726 (94%)	40 (5%)	6 (1%)	18	37

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	540	PHE
1	B	567[A]	GLU
1	B	567[B]	GLU
1	B	585	ASN
1	B	666	ARG
1	D	585	ASN

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	176/179 (98%)	153 (87%)	23 (13%)	4	10
1	B	180/179 (101%)	154 (86%)	26 (14%)	3	8
1	C	177/179 (99%)	159 (90%)	18 (10%)	7	18
1	D	175/179 (98%)	147 (84%)	28 (16%)	2	6
All	All	708/716 (99%)	613 (87%)	95 (13%)	4	9

All (95) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	526	GLN
1	A	539	SER
1	A	543	LEU
1	A	548	ARG
1	A	552	HIS
1	A	556	GLU
1	A	558	ARG
1	A	563	MET
1	A	564	PHE
1	A	567	GLU
1	A	582	ILE
1	A	589	LEU
1	A	608	LEU
1	A	612	ARG
1	A	627	THR
1	A	628	ILE
1	A	640	VAL
1	A	644	THR
1	A	649	GLU
1	A	662	CYS
1	A	666	ARG
1	A	674	ARG
1	A	701	ARG
1	B	521	ASP

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Mol	Chain	Res	Type
1	B	528	GLN
1	B	532	LYS
1	B	542	LYS
1	B	543	LEU
1	B	548	ARG
1	B	550	ARG
1	B	562	LYS
1	B	563	MET
1	B	564	PHE
1	B	582	ILE
1	B	587	ARG
1	B	589	LEU
1	B	608	LEU
1	B	630	LYS
1	B	636	GLN
1	B	640	VAL
1	B	645	LYS
1	B	647	ASN
1	B	648	LYS
1	B	662	CYS
1	B	665	THR
1	B	698	MET
1	B	705	THR
1	B	709	ASP
1	B	712	ASP
1	C	543	LEU
1	C	563	MET
1	C	582	ILE
1	C	590	VAL
1	C	592	SER
1	C	608	LEU
1	C	624[A]	ARG
1	C	624[B]	ARG
1	C	628	ILE
1	C	631	LYS
1	C	636	GLN
1	C	640	VAL
1	C	647	ASN
1	C	648	LYS
1	C	654	ASP
1	C	666	ARG
1	C	701	ARG

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Mol	Chain	Res	Type
1	C	712	ASP
1	D	535	GLU
1	D	543	LEU
1	D	556	GLU
1	D	562	LYS
1	D	564	PHE
1	D	566	GLU
1	D	567	GLU
1	D	582	ILE
1	D	589	LEU
1	D	593	MET
1	D	595	LEU
1	D	598	ASN
1	D	608	LEU
1	D	624	ARG
1	D	627	THR
1	D	636	GLN
1	D	640	VAL
1	D	645	LYS
1	D	648	LYS
1	D	649	GLU
1	D	656	SER
1	D	662	CYS
1	D	666	ARG
1	D	694	GLU
1	D	698	MET
1	D	701	ARG
1	D	709	ASP
1	D	712	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	536	GLN
1	A	585	ASN
1	A	602	ASN
1	B	528	GLN
1	B	549	GLN
1	C	526	GLN
1	D	602	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	CSO	D	586	1	3,6,7	1.07	0	1,6,8	0.06	0
1	CSO	A	586	1	3,6,7	1.26	0	1,6,8	0.36	0
1	CSO	B	586	1	3,6,7	0.66	0	1,6,8	0.33	0
1	CSO	C	586	1	3,6,7	0.81	0	1,6,8	0.30	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CSO	D	586	1	-	0/1/5/7	-
1	CSO	A	586	1	-	0/1/5/7	-
1	CSO	B	586	1	-	0/1/5/7	-
1	CSO	C	586	1	-	0/1/5/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	586	CSO	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

14 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	PCG	A	802	-	26,26,26	1.80	4 (15%)	39,41,41	2.09	9 (23%)
2	PCG	A	801	-	26,26,26	1.32	3 (11%)	39,41,41	2.60	11 (28%)
4	GOL	C	804	-	5,5,5	0.48	0	5,5,5	0.95	0
4	GOL	C	803	-	5,5,5	0.43	0	5,5,5	0.49	0
2	PCG	D	802	-	26,26,26	1.65	3 (11%)	39,41,41	2.18	9 (23%)
3	ACT	A	803	-	3,3,3	0.75	0	3,3,3	1.17	0
2	PCG	C	801	-	26,26,26	1.74	8 (30%)	39,41,41	2.00	8 (20%)
3	ACT	A	804	-	3,3,3	0.65	0	3,3,3	1.11	0
2	PCG	D	801	-	26,26,26	1.50	4 (15%)	39,41,41	2.23	9 (23%)
2	PCG	B	802	-	26,26,26	1.53	3 (11%)	39,41,41	2.32	12 (30%)
2	PCG	C	802	-	26,26,26	1.78	3 (11%)	39,41,41	2.17	11 (28%)
3	ACT	A	805	-	3,3,3	1.06	0	3,3,3	1.79	1 (33%)
2	PCG	B	801	-	26,26,26	1.48	5 (19%)	39,41,41	2.36	15 (38%)
3	ACT	C	805	-	3,3,3	0.74	0	3,3,3	1.36	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PCG	A	802	-	-	0/4/31/31	0/4/4/4
2	PCG	A	801	-	-	0/4/31/31	0/4/4/4
4	GOL	C	804	-	-	2/4/4/4	-
4	GOL	C	803	-	-	2/4/4/4	-
2	PCG	D	802	-	-	1/4/31/31	0/4/4/4
2	PCG	C	801	-	-	0/4/31/31	0/4/4/4
2	PCG	D	801	-	-	0/4/31/31	0/4/4/4
2	PCG	B	802	-	-	0/4/31/31	0/4/4/4
2	PCG	C	802	-	-	0/4/31/31	0/4/4/4
2	PCG	B	801	-	-	0/4/31/31	0/4/4/4

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	802	PCG	PA-O5'	5.94	1.64	1.57
2	C	802	PCG	PA-O5'	5.46	1.63	1.57
2	D	802	PCG	PA-O5'	4.92	1.63	1.57
2	C	802	PCG	C5-C4	3.91	1.49	1.38
2	C	801	PCG	PA-O5'	3.78	1.62	1.57
2	B	802	PCG	PA-O5'	3.62	1.61	1.57
2	C	802	PCG	PA-O3'	3.59	1.63	1.57
2	B	802	PCG	C5-C4	3.58	1.48	1.38
2	A	802	PCG	C5-C4	3.57	1.48	1.38
2	D	802	PCG	C5-C4	3.43	1.48	1.38
2	D	801	PCG	C5-C4	3.42	1.48	1.38
2	B	801	PCG	PA-O5'	3.34	1.61	1.57
2	A	801	PCG	C5-N7	-3.30	1.32	1.39
2	C	801	PCG	O5'-C5'	-3.26	1.41	1.46
2	C	801	PCG	C5-C4	3.09	1.47	1.38
2	B	801	PCG	C6-N1	-3.07	1.33	1.38
2	A	802	PCG	PA-O3'	2.84	1.62	1.57
2	D	802	PCG	PA-O3'	2.82	1.62	1.57
2	D	801	PCG	PA-O3'	2.71	1.62	1.57
2	D	801	PCG	PA-O5'	2.69	1.60	1.57
2	C	801	PCG	O3'-C3'	-2.63	1.40	1.44
2	A	801	PCG	C5-C4	2.56	1.45	1.38
2	B	802	PCG	PA-O3'	2.54	1.62	1.57
2	C	801	PCG	C6-N1	-2.53	1.34	1.38
2	C	801	PCG	PA-O3'	2.48	1.61	1.57
2	B	801	PCG	C5-N7	-2.44	1.34	1.39
2	B	801	PCG	C5-C4	2.33	1.45	1.38
2	B	801	PCG	O3'-C3'	-2.24	1.41	1.44
2	C	801	PCG	C5-N7	-2.23	1.34	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	PCG	C6-N1	-2.22	1.34	1.38
2	C	801	PCG	C4-N9	-2.21	1.32	1.38
2	A	802	PCG	C6-N1	-2.10	1.34	1.38
2	D	801	PCG	O6-C6	2.04	1.27	1.23

All (85) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	PCG	O3'-C3'-C4'	-8.11	104.59	110.71
2	B	801	PCG	C5-C4-N3	-6.73	117.68	128.39
2	D	802	PCG	C5-C4-N3	-6.65	117.80	128.39
2	B	802	PCG	C5-C4-N3	-6.26	118.43	128.39
2	C	801	PCG	O3'-C3'-C4'	-6.23	106.01	110.71
2	A	801	PCG	C5-C4-N3	-6.15	118.61	128.39
2	C	802	PCG	C5-C4-N3	-6.04	118.77	128.39
2	B	802	PCG	O3'-C3'-C4'	-5.96	106.21	110.71
2	A	802	PCG	C5-C4-N3	-5.93	118.94	128.39
2	D	801	PCG	C5-C4-N3	-5.85	119.08	128.39
2	A	801	PCG	C2-N3-C4	5.67	122.06	112.30
2	D	802	PCG	C2-N3-C4	5.62	121.98	112.30
2	B	801	PCG	C2-N3-C4	5.46	121.71	112.30
2	D	801	PCG	O3'-C3'-C4'	-5.37	106.66	110.71
2	B	802	PCG	C2-N3-C4	5.34	121.50	112.30
2	C	802	PCG	C2-N3-C4	5.32	121.47	112.30
2	A	802	PCG	N9-C4-N3	5.17	136.29	125.95
2	C	801	PCG	C5-C4-N3	-5.16	120.18	128.39
2	A	802	PCG	C2-N3-C4	5.14	121.16	112.30
2	A	801	PCG	N9-C4-N3	4.96	135.87	125.95
2	D	801	PCG	O1A-PA-O2A	4.94	123.73	108.56
2	D	802	PCG	N9-C4-N3	4.87	135.69	125.95
2	D	801	PCG	C2-N3-C4	4.86	120.67	112.30
2	B	801	PCG	N9-C4-N3	4.78	135.51	125.95
2	B	801	PCG	O3'-C3'-C4'	-4.66	107.19	110.71
2	C	802	PCG	N9-C4-N3	4.61	135.18	125.95
2	B	802	PCG	N9-C4-N3	4.54	135.02	125.95
2	D	801	PCG	N9-C4-N3	4.37	134.70	125.95
2	C	802	PCG	O1A-PA-O2A	4.32	121.83	108.56
2	D	802	PCG	O1A-PA-O2A	4.08	121.09	108.56
2	A	801	PCG	O1A-PA-O2A	4.05	121.01	108.56
2	C	801	PCG	C2-N3-C4	4.04	119.26	112.30
2	A	802	PCG	O1A-PA-O2A	3.97	120.77	108.56
2	A	801	PCG	O5'-PA-O3'	-3.84	100.56	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	PCG	O1A-PA-O2A	3.78	120.16	108.56
2	A	801	PCG	O6-C6-C5	-3.77	116.59	126.53
2	D	801	PCG	O3'-PA-O2A	-3.72	102.41	110.39
2	B	802	PCG	O1A-PA-O2A	3.66	119.79	108.56
2	C	801	PCG	O1A-PA-O2A	3.52	119.39	108.56
2	C	801	PCG	N9-C4-N3	3.46	132.88	125.95
2	D	801	PCG	C6-C5-N7	3.31	136.31	130.29
2	C	802	PCG	C6-C5-N7	3.30	136.29	130.29
2	B	802	PCG	O3'-C3'-C2'	3.27	118.81	115.61
2	D	802	PCG	O4'-C1'-N9	3.13	115.45	108.36
2	D	802	PCG	C6-C5-N7	3.13	135.98	130.29
2	A	802	PCG	C6-C5-N7	3.07	135.88	130.29
2	C	802	PCG	O3'-C3'-C2'	2.86	118.41	115.61
2	A	802	PCG	O6-C6-C5	-2.81	119.11	126.53
2	B	801	PCG	C4'-O4'-C1'	-2.76	103.37	109.47
2	B	802	PCG	C6-C5-N7	2.69	135.19	130.29
2	D	801	PCG	C4-C5-N7	-2.68	106.42	110.67
2	C	802	PCG	O3'-C3'-C4'	2.63	112.69	110.71
2	B	801	PCG	C4-C5-N7	-2.62	106.52	110.67
2	A	801	PCG	C5-C6-N1	2.60	119.86	113.25
2	B	801	PCG	O6-C6-C5	-2.58	119.73	126.53
2	D	802	PCG	C4-C5-N7	-2.56	106.61	110.67
2	B	801	PCG	C6-C5-N7	2.51	134.87	130.29
2	B	802	PCG	O6-C6-C5	-2.51	119.91	126.53
2	C	802	PCG	C4-C5-N7	-2.49	106.73	110.67
2	C	801	PCG	C6-C5-N7	2.47	134.78	130.29
2	C	801	PCG	O6-C6-C5	-2.46	120.03	126.53
2	B	802	PCG	C5'-C4'-C3'	-2.46	107.94	112.56
2	B	801	PCG	O5'-PA-O3'	-2.44	102.44	105.70
3	A	805	ACT	OXT-C-CH3	2.41	125.17	115.05
2	D	802	PCG	O5'-PA-O2A	-2.36	105.00	110.44
2	A	802	PCG	C5-C6-N1	2.34	119.21	113.25
2	C	802	PCG	O6-C6-C5	-2.33	120.39	126.53
2	C	801	PCG	O5'-PA-O2A	-2.32	105.08	110.44
2	B	801	PCG	O4'-C4'-C5'	2.31	119.36	112.38
2	A	801	PCG	O3'-C3'-C2'	2.29	117.86	115.61
2	B	802	PCG	O5'-PA-O2A	-2.22	105.31	110.44
2	A	801	PCG	C6-C5-N7	2.20	134.30	130.29
2	A	801	PCG	O5'-PA-O2A	-2.20	105.37	110.44
2	B	801	PCG	C2-N1-C6	-2.17	121.17	125.11
2	B	802	PCG	C4-C5-N7	-2.16	107.25	110.67
2	A	802	PCG	O2'-C2'-C3'	-2.15	105.17	111.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	PCG	O5'-PA-O2A	-2.13	105.53	110.44
2	B	802	PCG	O5'-PA-O3'	-2.13	102.86	105.70
2	B	801	PCG	O3'-C3'-C2'	2.13	117.69	115.61
2	C	802	PCG	C5'-C4'-C3'	-2.06	108.69	112.56
2	C	802	PCG	C2'-C1'-N9	-2.06	107.52	113.25
2	D	802	PCG	O6-C6-C5	-2.05	121.12	126.53
2	B	801	PCG	C8-N7-C5	2.04	107.90	104.26
2	D	801	PCG	O4'-C4'-C3'	-2.03	100.63	104.92
2	A	802	PCG	O3'-C3'-C2'	-2.01	113.63	115.61

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	803	GOL	C1-C2-C3-O3
4	C	804	GOL	C1-C2-C3-O3
4	C	804	GOL	O2-C2-C3-O3
4	C	803	GOL	O2-C2-C3-O3
2	D	802	PCG	C2'-C1'-N9-C4

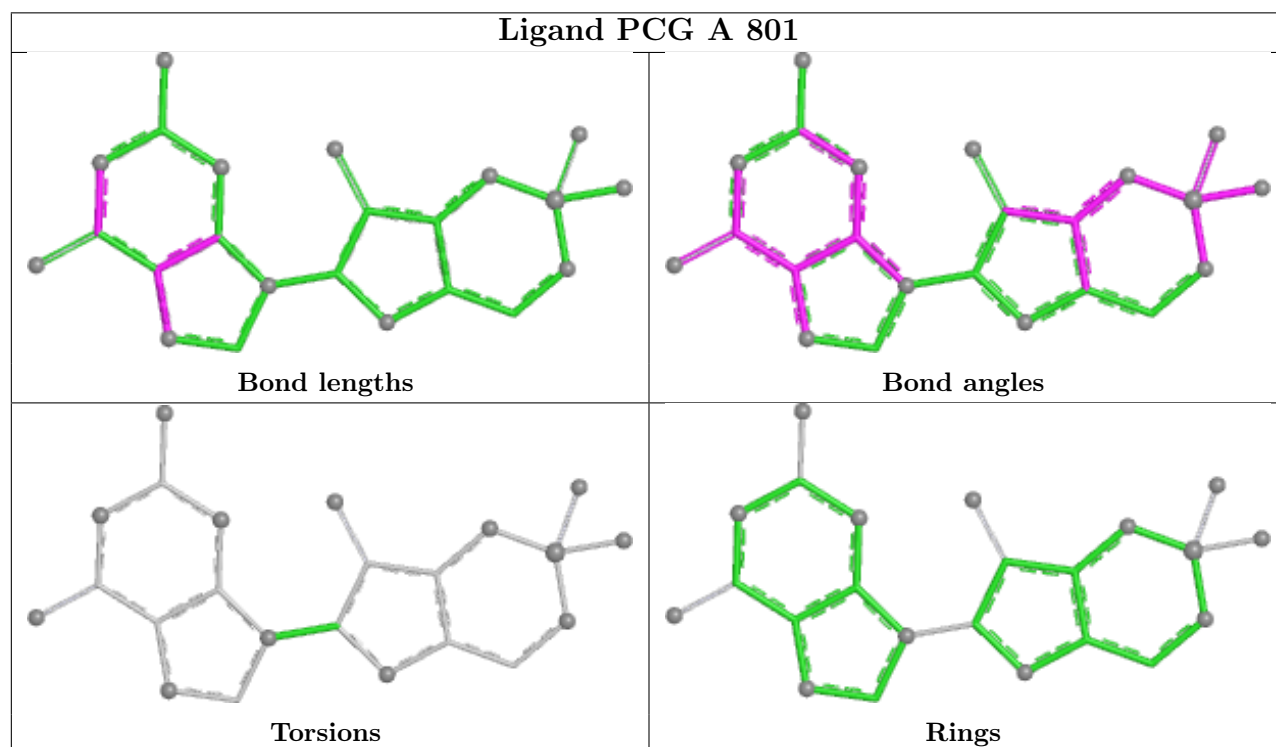
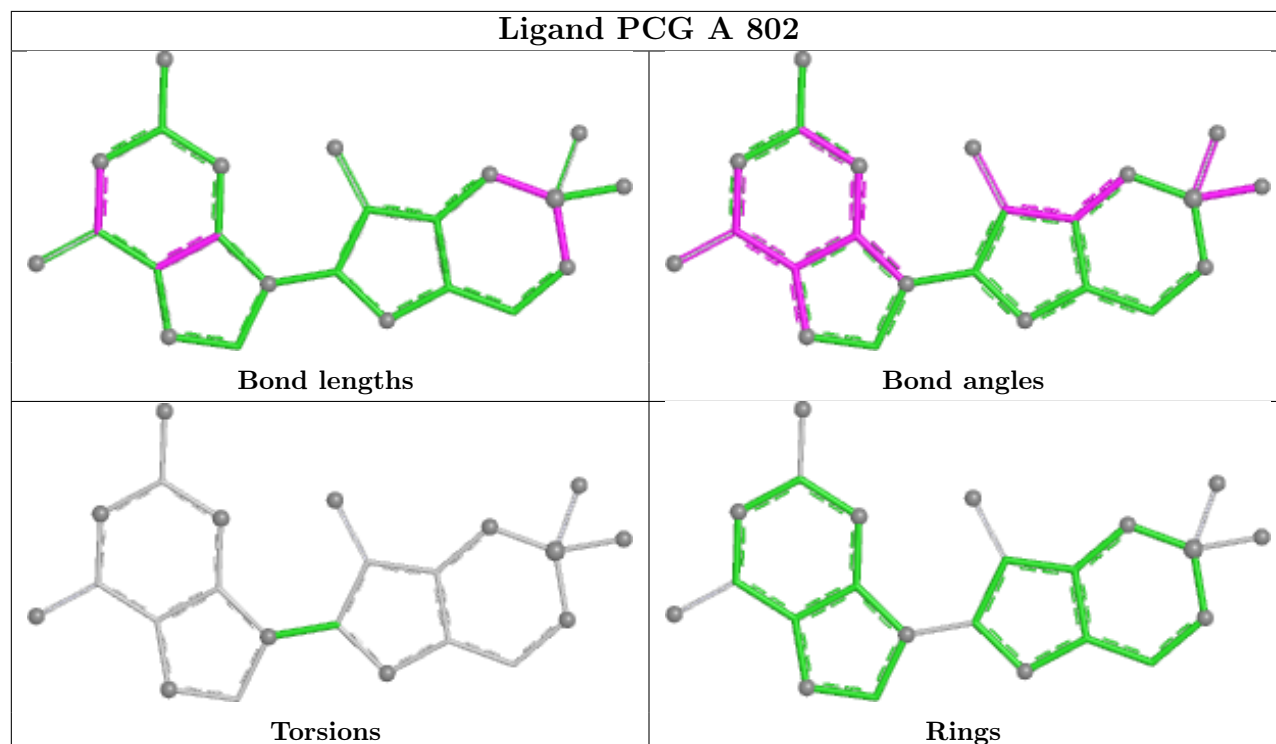
There are no ring outliers.

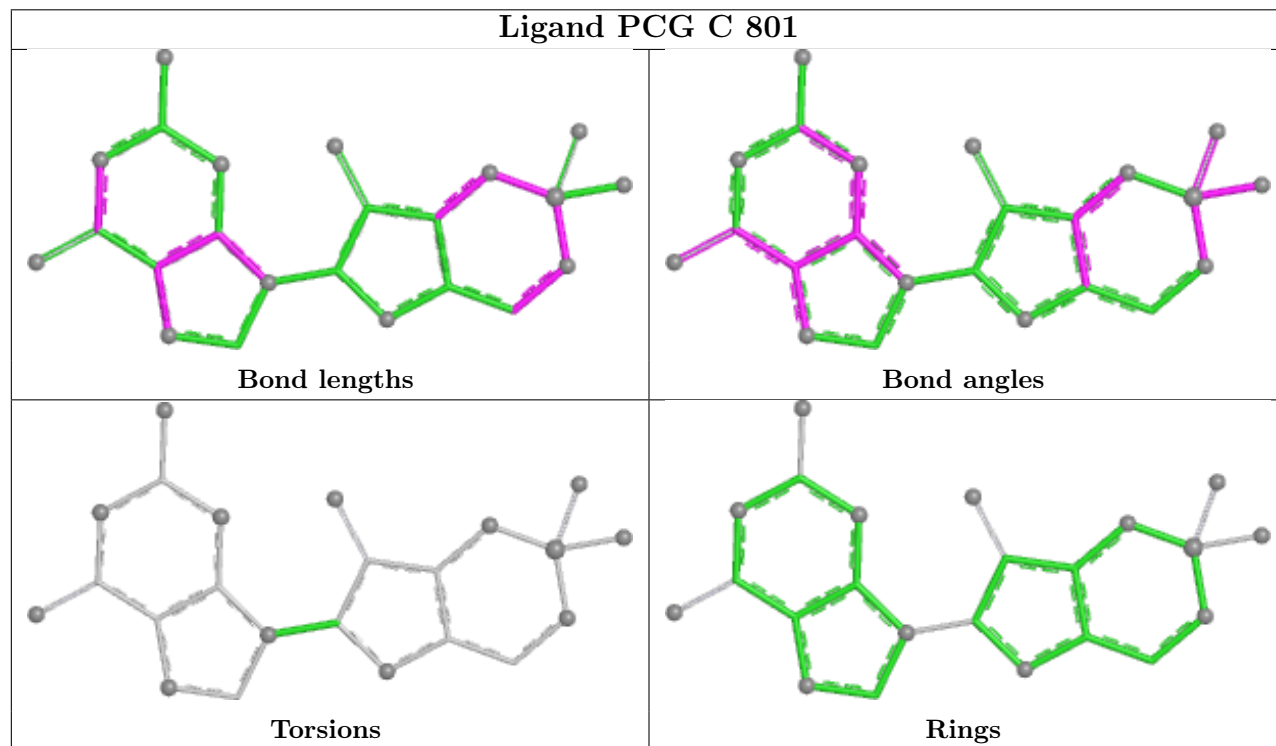
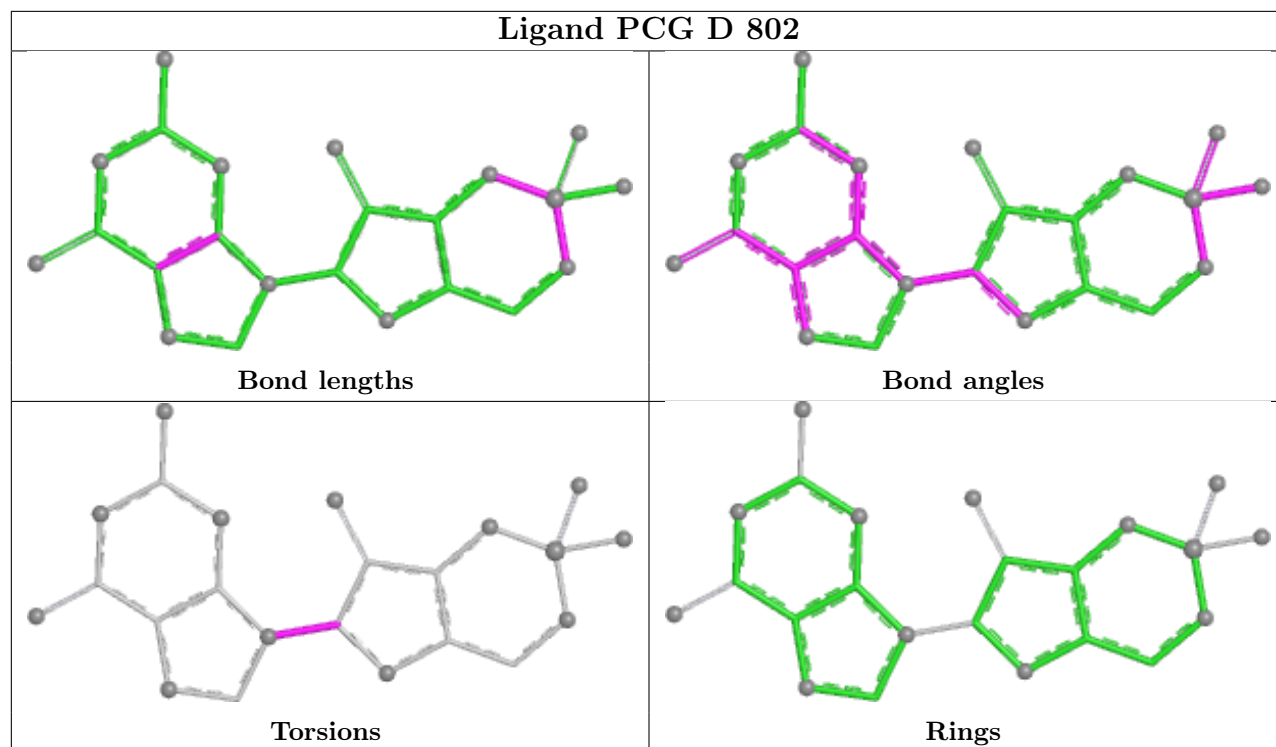
8 monomers are involved in 13 short contacts:

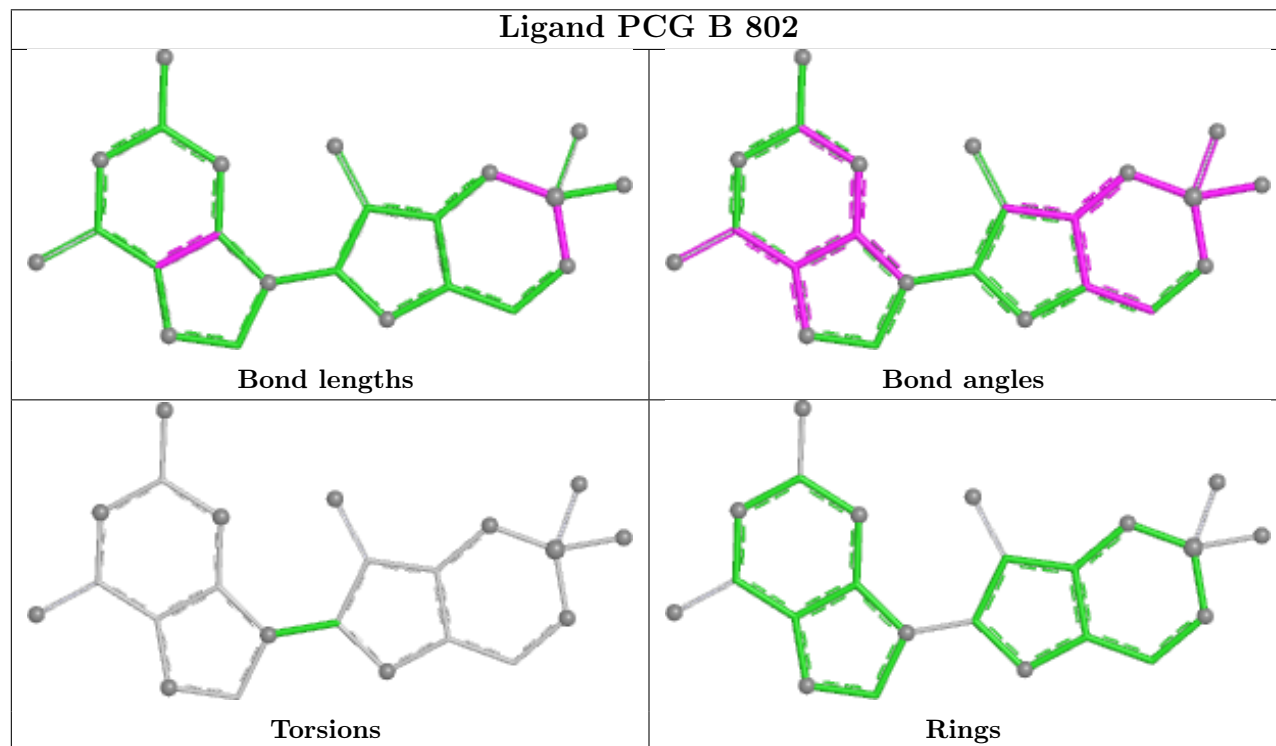
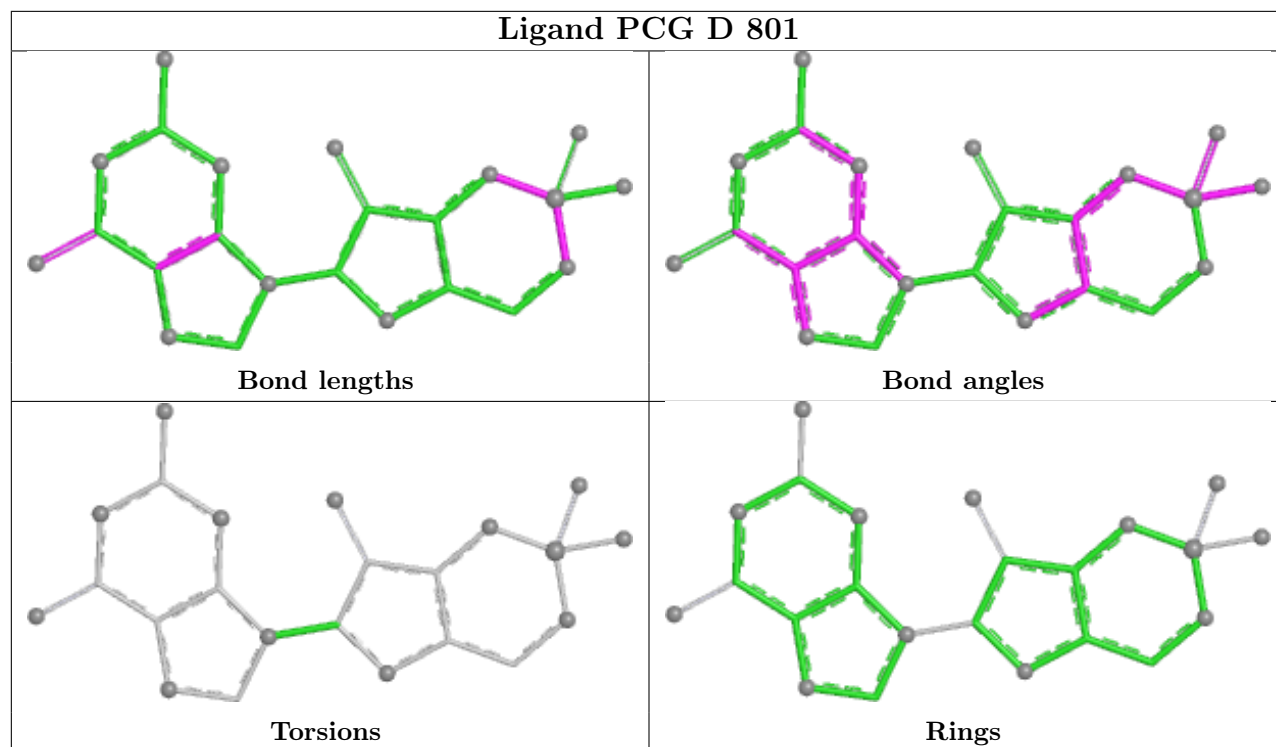
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	PCG	1	0
4	C	804	GOL	3	0
2	D	802	PCG	1	0
3	A	803	ACT	2	0
2	C	801	PCG	1	0
3	A	804	ACT	3	0
2	B	802	PCG	1	0
2	B	801	PCG	1	0

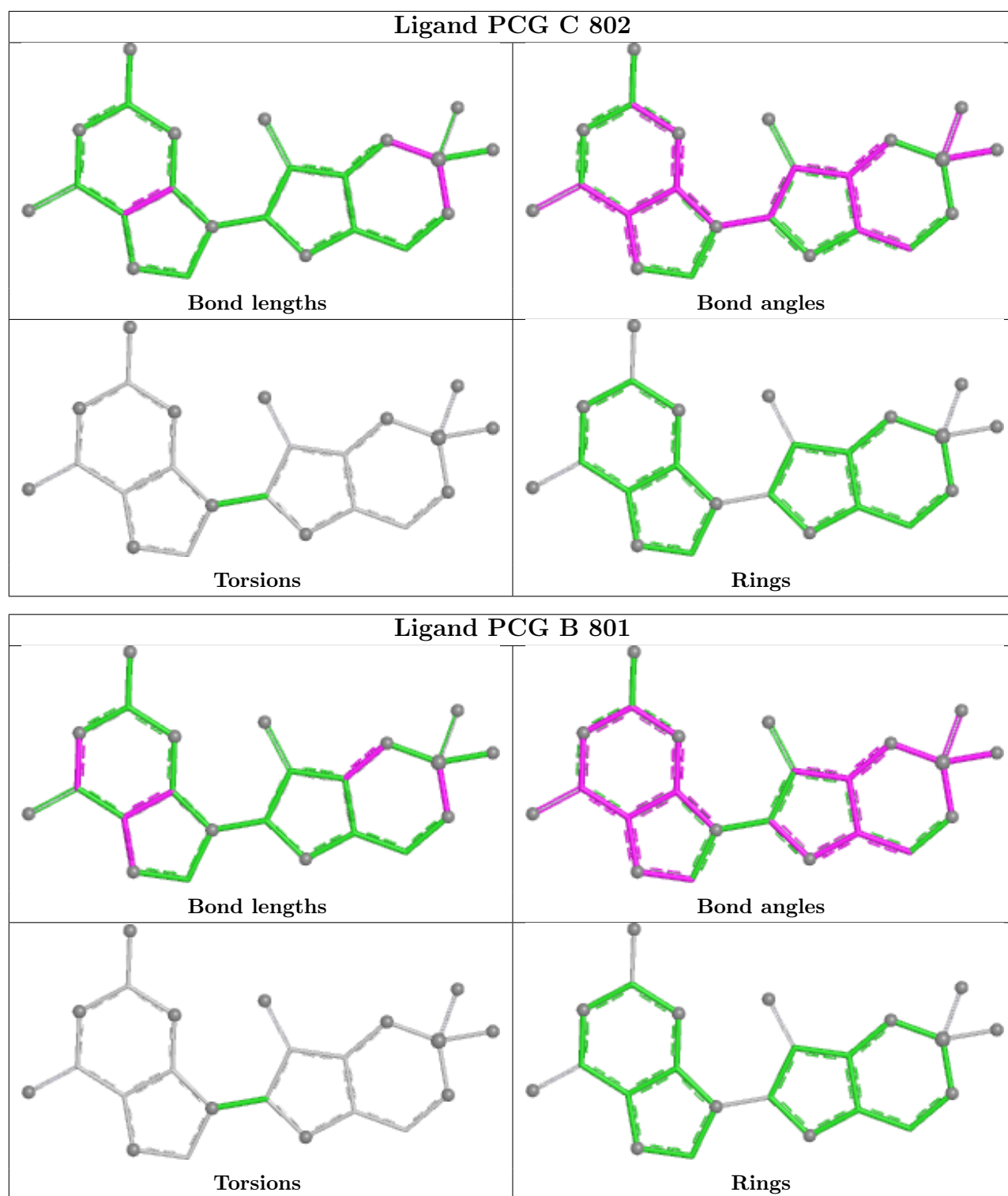
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	192/199 (96%)	-0.25	3 (1%) 70 68	12, 29, 61, 75	2 (1%)
1	B	192/199 (96%)	-0.10	7 (3%) 46 42	14, 34, 65, 85	6 (3%)
1	C	192/199 (96%)	-0.33	2 (1%) 79 78	11, 31, 58, 72	3 (1%)
1	D	192/199 (96%)	0.12	6 (3%) 51 48	22, 41, 67, 78	1 (0%)
All	All	768/796 (96%)	-0.14	18 (2%) 61 58	11, 34, 65, 85	12 (1%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	521	ASP	3.3
1	C	521	ASP	3.2
1	B	521	ASP	3.1
1	B	567[A]	GLU	3.1
1	A	698[A]	MET	3.1
1	D	598	ASN	2.7
1	D	593	MET	2.5
1	B	645	LYS	2.4
1	C	713	ARG	2.4
1	B	524	ARG	2.4
1	B	708	LEU	2.4
1	D	648	LYS	2.3
1	D	666	ARG	2.3
1	B	713	ARG	2.2
1	B	698	MET	2.2
1	A	713	ARG	2.2
1	D	698	MET	2.2
1	A	564	PHE	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	CSO	D	586	7/8	0.88	0.12	45,46,46,48	0
1	CSO	C	586	7/8	0.94	0.10	16,32,33,33	0
1	CSO	B	586	7/8	0.94	0.09	32,36,37,37	0
1	CSO	A	586	7/8	0.95	0.08	29,31,32,34	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

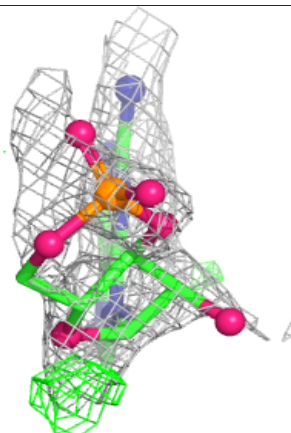
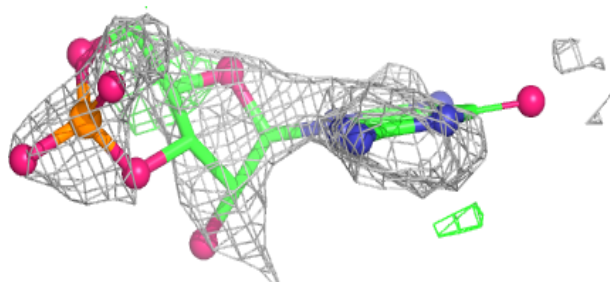
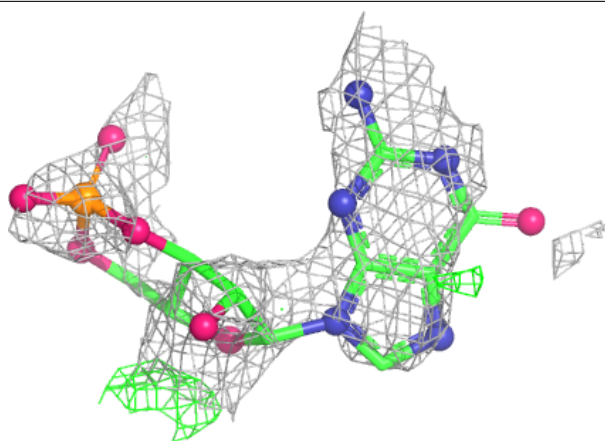
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	PCG	D	802	23/23	0.72	0.27	82,84,86,86	23
4	GOL	C	803	6/6	0.80	0.18	63,65,66,66	0
4	GOL	C	804	6/6	0.81	0.17	61,62,63,64	0
2	PCG	B	802	23/23	0.82	0.18	62,64,65,65	23
3	ACT	A	803	4/4	0.85	0.20	68,68,68,69	0
2	PCG	A	802	23/23	0.88	0.10	51,56,59,60	0
2	PCG	C	802	23/23	0.88	0.11	59,63,65,65	0
3	ACT	A	805	4/4	0.92	0.11	27,28,28,28	0
3	ACT	C	805	4/4	0.92	0.14	65,66,66,66	0
3	ACT	A	804	4/4	0.94	0.11	31,31,32,33	0
2	PCG	B	801	23/23	0.97	0.07	17,24,29,32	0
2	PCG	A	801	23/23	0.98	0.05	12,17,20,21	0
2	PCG	D	801	23/23	0.98	0.05	22,28,38,38	0
2	PCG	C	801	23/23	0.98	0.05	21,25,28,30	0

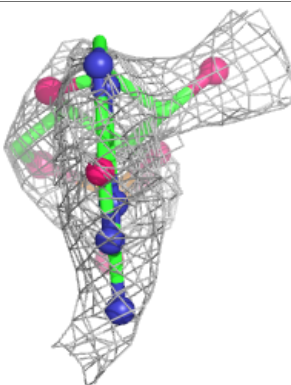
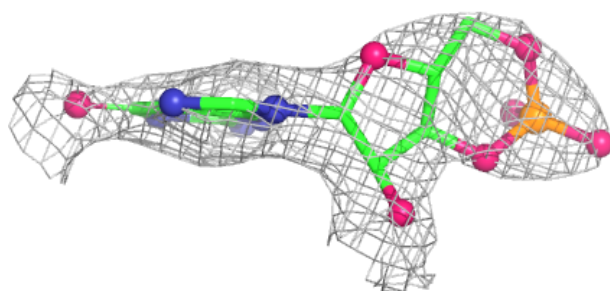
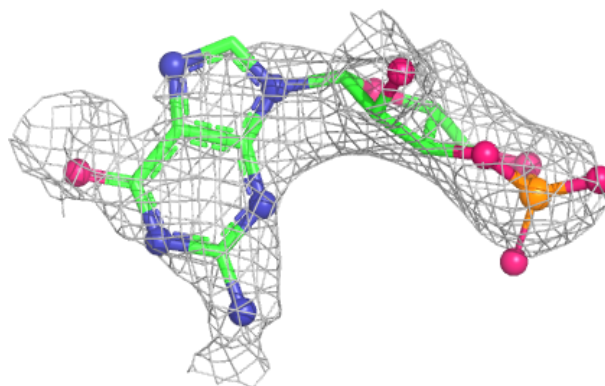
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around PCG D 802:

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

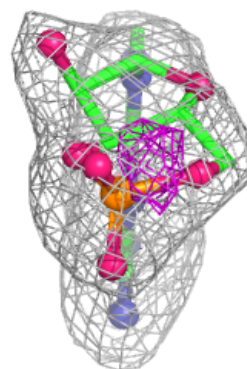
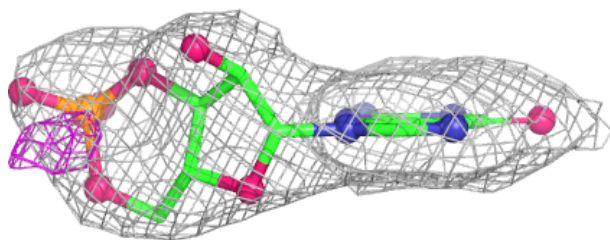
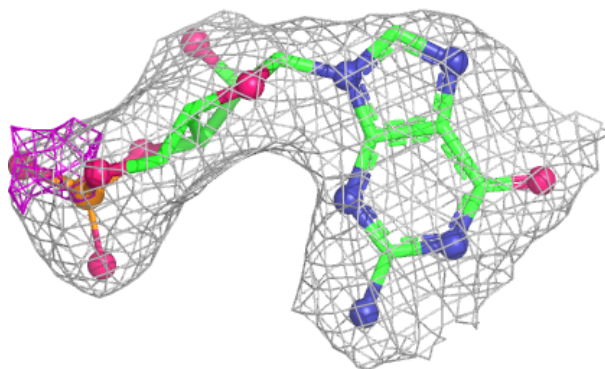
**Electron density around PCG B 802:**

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

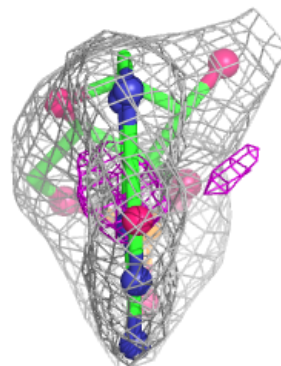
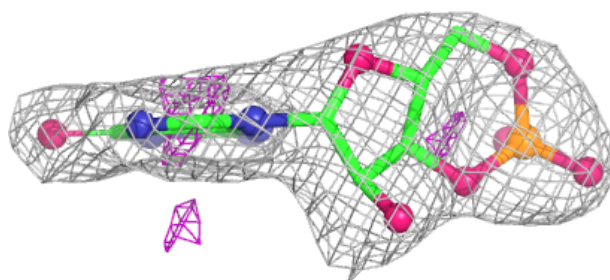
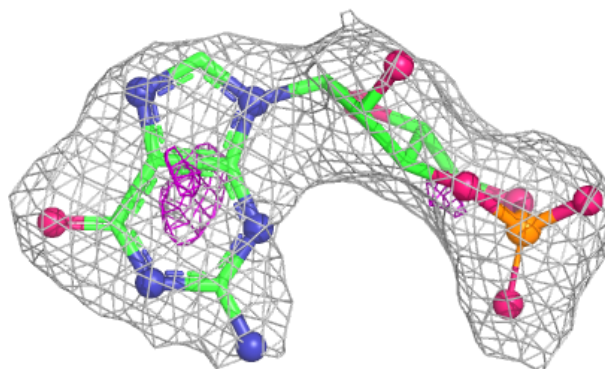


Electron density around PCG A 802:

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

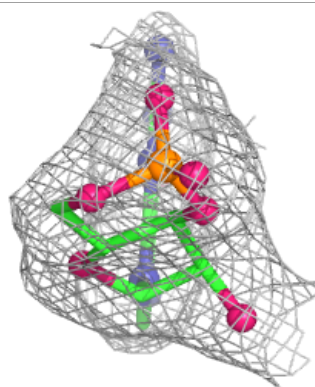
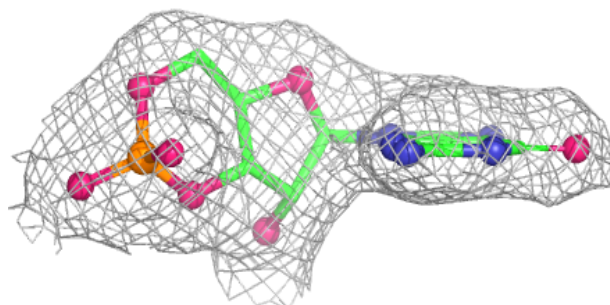
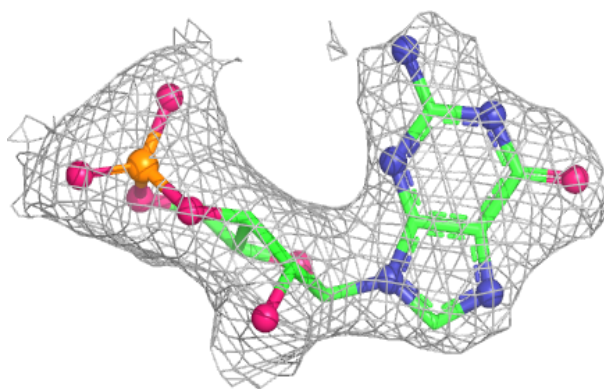
**Electron density around PCG C 802:**

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

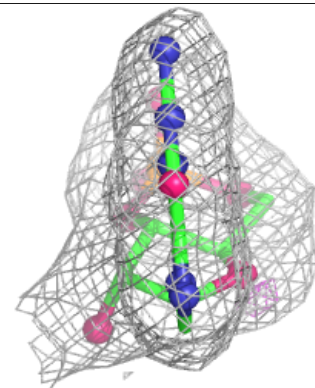
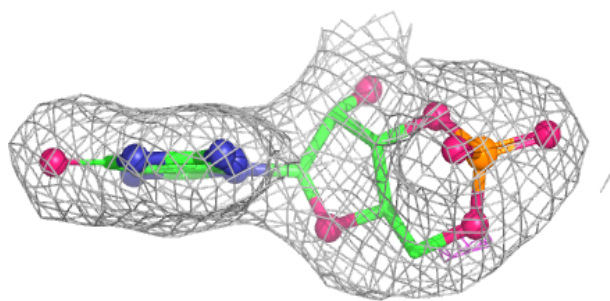
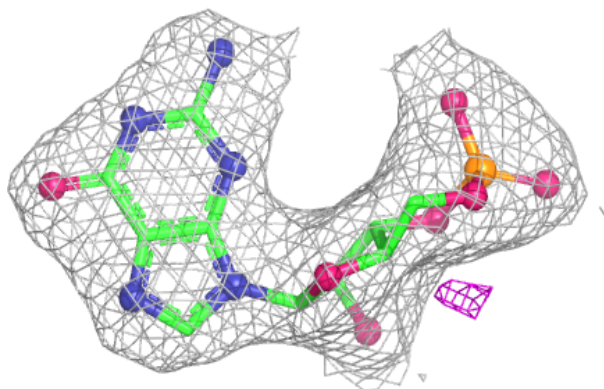


Electron density around PCG B 801:

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

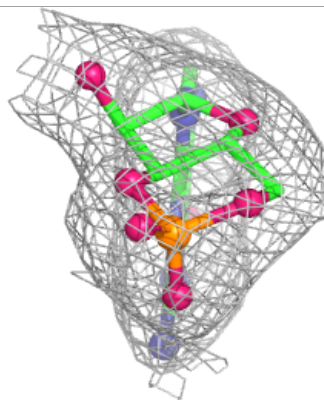
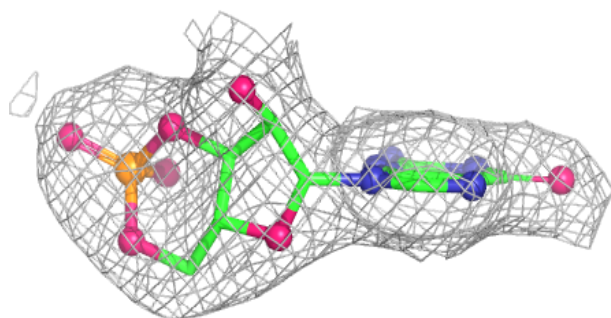
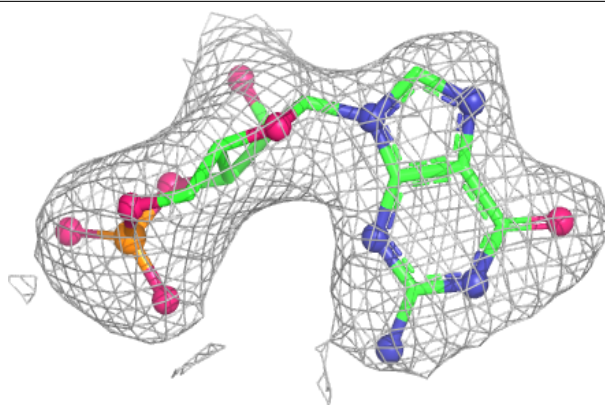
**Electron density around PCG A 801:**

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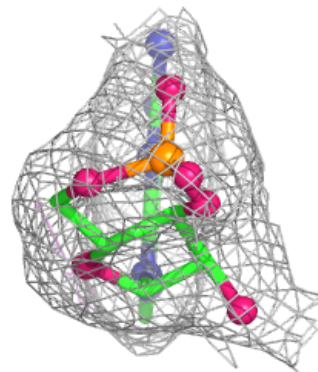
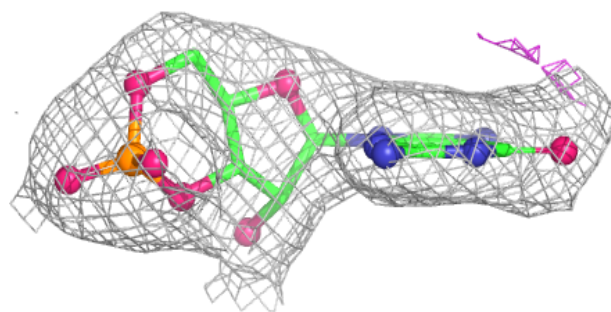
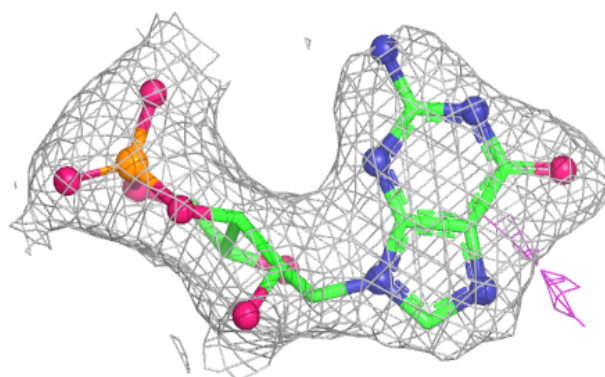


Electron density around PCG D 801:

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and green (positive)

**Electron density around PCG C 801:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.