



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 12, 2022 – 10:08 AM JST

PDB ID : 8GXF  
Title : Pseudomonas flexibilis GCN5 family acetyltransferase  
Authors : Song, Y.J.; Bao, R.  
Deposited on : 2022-09-20  
Resolution : 3.04 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](mailto: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 : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.31.2  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.2

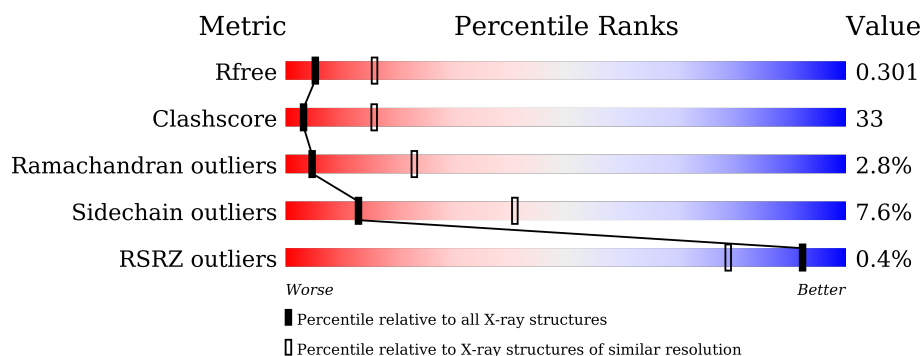
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.04 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))
$R_{free}$	130704	2752 (3.08-3.00)
Clashscore	141614	3096 (3.08-3.00)
Ramachandran outliers	138981	2986 (3.08-3.00)
Sidechain outliers	138945	2988 (3.08-3.00)
RSRZ outliers	127900	2636 (3.08-3.00)

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  $\geq 3$ , 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	193	<div> <div></div> <div>46%</div> <div>44%</div> <div>7%</div> <div>• •</div> </div>
1	B	193	<div> <div></div> <div>49%</div> <div>38%</div> <div>6%</div> <div>• •</div> </div>
1	C	193	<div> <div>%</div> <div>46%</div> <div>41%</div> <div>9%</div> <div>• •</div> </div>
1	D	193	<div> <div>%</div> <div>54%</div> <div>35%</div> <div>6%</div> <div>• •</div> </div>

## 2 Entry composition [i](#)

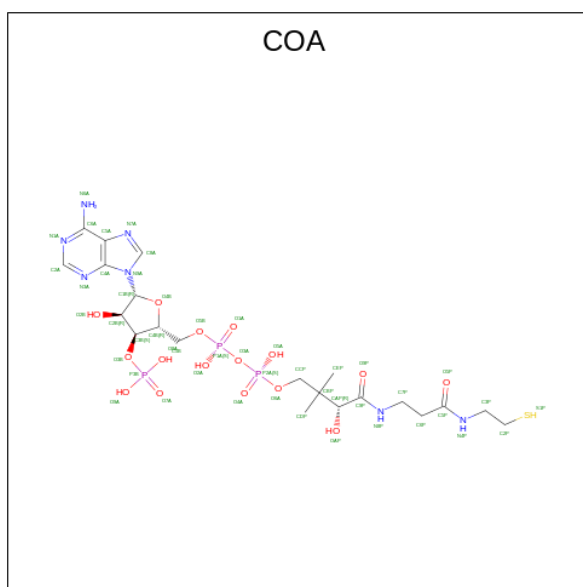
There are 3 unique types of molecules in this entry. The entry contains 6245 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GCN5 family acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	187	Total	C	N	O	S	0	0	0
			1504	944	283	272	5			
1	A	187	Total	C	N	O	S	0	0	0
			1504	944	283	272	5			
1	C	187	Total	C	N	O	S	0	0	0
			1504	944	283	272	5			
1	D	187	Total	C	N	O	S	0	0	0
			1504	944	283	272	5			

- Molecule 2 is COENZYME A (three-letter code: COA) (formula:  $C_{21}H_{36}N_7O_{16}P_3S$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P S	0	0
			48	21	7	16	3 1		
2	A	1	Total	C	N	O	P S	0	0
			48	21	7	16	3 1		

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
2	D	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

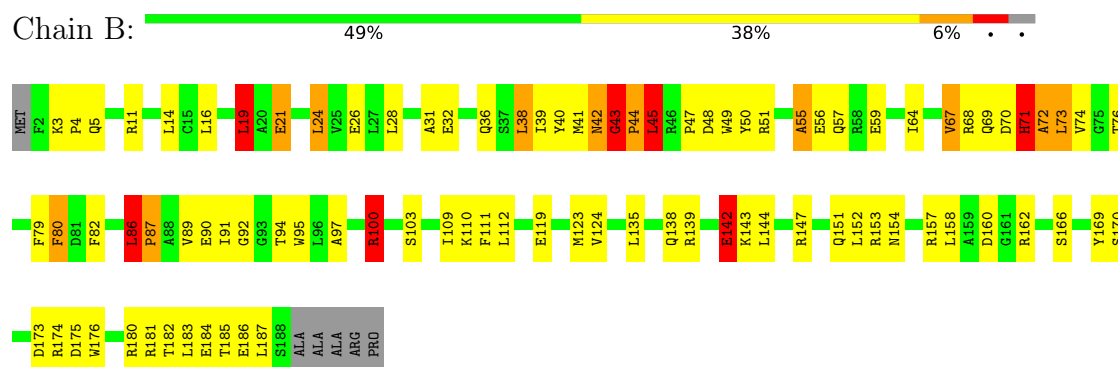
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	6	Total	O	0	0
			6	6		
3	A	13	Total	O	0	0
			13	13		
3	C	11	Total	O	0	0
			11	11		
3	D	7	Total	O	0	0
			7	7		

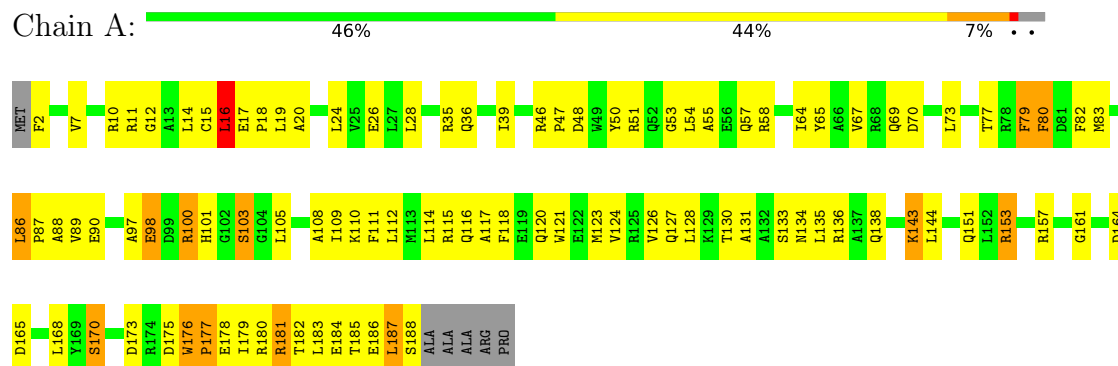
### 3 Residue-property plots

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.

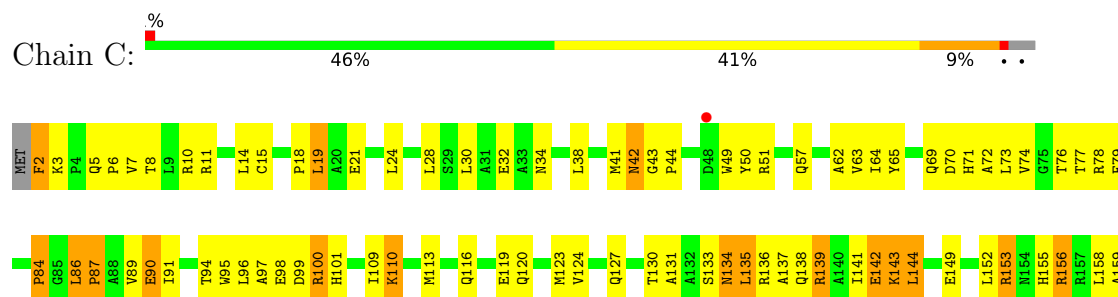
#### • Molecule 1: GCN5 family acetyltransferase

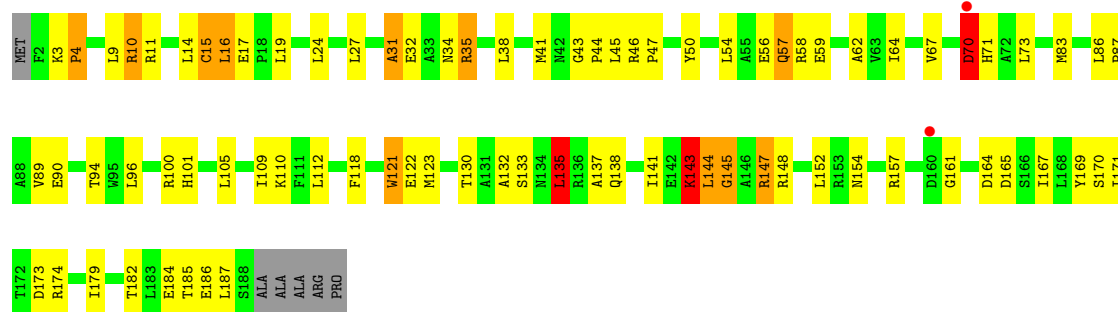


#### • Molecule 1: GCN5 family acetyltransferase



#### • Molecule 1: GCN5 family acetyltransferase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	43.89Å 99.01Å 98.82Å 90.00° 101.87° 90.00°	Depositor
Resolution (Å)	42.95 – 3.04 49.51 – 3.04	Depositor EDS
% Data completeness (in resolution range)	99.0 (42.95-3.04) 92.6 (49.51-3.04)	Depositor EDS
$R_{merge}$	0.47	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.48 (at 3.07Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, $R_{free}$	0.225 , 0.302 0.225 , 0.301	Depositor DCC
$R_{free}$ test set	1585 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.4	Xtriage
Anisotropy	0.573	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 55.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.042 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.85	EDS
Total number of atoms	6245	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.13% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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

## 5 Model quality

### 5.1 Standard geometry

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

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	1.08	4/1532 (0.3%)	0.95	2/2070 (0.1%)
1	B	0.99	7/1532 (0.5%)	1.06	14/2070 (0.7%)
1	C	0.98	3/1532 (0.2%)	0.96	5/2070 (0.2%)
1	D	0.72	3/1532 (0.2%)	0.81	6/2070 (0.3%)
All	All	0.95	17/6128 (0.3%)	0.95	27/8280 (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	D	0	1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	90	GLU	CD-OE1	-8.10	1.16	1.25
1	A	79	PHE	CG-CD1	-6.29	1.29	1.38
1	B	21	GLU	CG-CD	6.24	1.61	1.51
1	B	40	TYR	CG-CD1	-5.75	1.31	1.39
1	B	142	GLU	CD-OE1	-5.68	1.19	1.25
1	A	20	ALA	C-O	-5.55	1.12	1.23
1	B	142	GLU	CD-OE2	-5.49	1.19	1.25
1	C	184	GLU	CD-OE1	-5.48	1.19	1.25
1	D	17	GLU	CD-OE1	-5.38	1.19	1.25
1	D	17	GLU	CD-OE2	-5.32	1.19	1.25
1	A	17	GLU	CD-OE1	-5.29	1.19	1.25
1	B	40	TYR	CG-CD2	-5.24	1.32	1.39
1	B	87	PRO	C-O	-5.22	1.12	1.23
1	D	145	GLY	C-O	-5.15	1.15	1.23

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	67	VAL	C-O	-5.13	1.13	1.23
1	C	142	GLU	CD-OE1	-5.12	1.20	1.25
1	A	176	TRP	CE3-CZ3	-5.07	1.29	1.38

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	100	ARG	NE-CZ-NH2	9.71	125.16	120.30
1	A	16	LEU	CB-CG-CD2	-8.61	96.36	111.00
1	B	71	HIS	N-CA-C	-8.12	89.08	111.00
1	B	87	PRO	CA-N-CD	-6.03	103.06	111.50
1	A	79	PHE	CB-CA-C	-5.93	98.54	110.40
1	B	72	ALA	N-CA-C	-5.91	95.03	111.00
1	B	100	ARG	NE-CZ-NH1	-5.91	117.34	120.30
1	D	144	LEU	CB-CG-CD2	-5.91	100.96	111.00
1	C	187	LEU	N-CA-C	-5.80	95.33	111.00
1	D	135	LEU	CA-CB-CG	5.79	128.61	115.30
1	C	87	PRO	CA-N-CD	-5.67	103.56	111.50
1	B	100	ARG	CD-NE-CZ	5.56	131.38	123.60
1	D	143	LYS	CD-CE-NZ	-5.52	99.01	111.70
1	B	21	GLU	OE1-CD-OE2	-5.51	116.68	123.30
1	C	19	LEU	CA-CB-CG	5.45	127.84	115.30
1	B	43	GLY	N-CA-C	-5.34	99.76	113.10
1	C	84	PRO	N-CA-C	5.27	125.80	112.10
1	B	73	LEU	N-CA-C	-5.23	96.89	111.00
1	C	89	VAL	CB-CA-C	-5.22	101.49	111.40
1	B	38	LEU	CB-CG-CD1	-5.21	102.14	111.00
1	D	70	ASP	CB-CG-OD2	5.19	122.97	118.30
1	D	15	CYS	CA-CB-SG	5.18	123.32	114.00
1	B	74	VAL	CB-CA-C	5.17	121.22	111.40
1	B	19	LEU	CA-CB-CG	5.16	127.17	115.30
1	B	24	LEU	CA-CB-CG	-5.11	103.55	115.30
1	B	45	LEU	CB-CG-CD2	-5.09	102.35	111.00
1	D	16	LEU	CA-CB-CG	5.01	126.82	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	135	LEU	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	1504	0	1509	111	1
1	B	1504	0	1509	120	1
1	C	1504	0	1509	111	0
1	D	1504	0	1509	78	0
2	A	48	0	32	7	0
2	B	48	0	32	5	0
2	C	48	0	32	11	0
2	D	48	0	32	11	0
3	A	13	0	0	0	0
3	B	6	0	0	0	0
3	C	11	0	0	0	0
3	D	7	0	0	0	0
All	All	6245	0	6164	413	1

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 33.

All (413) 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:24:LEU:HD21	1:B:50:TYR:CB	1.16	1.59
1:B:24:LEU:CD2	1:B:50:TYR:HB2	1.35	1.51
1:A:175:ASP:O	1:A:179:ILE:HD13	1.25	1.28
1:B:24:LEU:CD2	1:B:50:TYR:CB	2.02	1.22
2:D:201:COA:O4B	2:D:201:COA:C1B	1.63	1.19
2:B:201:COA:O4B	2:B:201:COA:C1B	1.63	1.18
1:B:24:LEU:CD2	1:B:50:TYR:CG	2.28	1.17
1:A:175:ASP:O	1:A:179:ILE:CD1	1.97	1.12
1:B:26:GLU:OE1	1:B:71:HIS:HE1	1.34	1.11
1:A:14:LEU:HD13	1:A:105:LEU:HD11	1.26	1.07
1:B:97:ALA:HB3	1:B:100:ARG:NH2	1.69	1.05
1:A:182:THR:O	1:A:185:THR:OG1	1.75	1.05
1:A:14:LEU:CD1	1:A:105:LEU:HD11	1.88	1.03
1:C:24:LEU:HD12	1:C:50:TYR:CB	1.90	1.02
1:C:41:MET:O	1:C:159:ALA:HB2	1.61	0.99

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:LEU:HD21	1:B:50:TYR:CG	1.93	0.99
1:B:26:GLU:OE1	1:B:71:HIS:CE1	2.16	0.98
1:A:116:GLN:OE1	1:A:120:GLN:OE1	1.82	0.98
1:B:24:LEU:HD21	1:B:50:TYR:HB3	1.46	0.97
1:A:116:GLN:OE1	1:A:120:GLN:CD	2.03	0.97
1:B:24:LEU:CD2	1:B:50:TYR:CD1	2.51	0.94
1:D:35:ARG:HE	1:D:45:LEU:HD13	1.31	0.93
1:B:68:ARG:HG3	1:B:73:LEU:HD23	1.52	0.92
1:A:14:LEU:HD13	1:A:105:LEU:CD1	1.99	0.92
1:B:24:LEU:HD23	1:B:50:TYR:CG	2.02	0.91
1:A:185:THR:HG1	1:A:186:GLU:H	1.16	0.90
1:B:89:VAL:HG22	1:B:123:MET:HE1	1.53	0.90
1:B:138:GLN:O	1:B:142:GLU:HG3	1.72	0.90
1:B:42:ASN:O	1:B:42:ASN:ND2	2.07	0.88
1:A:116:GLN:OE1	1:A:120:GLN:NE2	2.07	0.88
1:C:182:THR:O	1:C:185:THR:OG1	1.92	0.87
1:B:97:ALA:HB3	1:B:100:ARG:HH21	1.35	0.86
1:C:24:LEU:CD1	1:C:50:TYR:HB3	2.04	0.86
1:B:24:LEU:HD23	1:B:50:TYR:CD1	2.12	0.84
1:C:139:ARG:NH1	1:C:143:LYS:HG2	1.91	0.84
1:C:110:LYS:NZ	1:C:110:LYS:HB3	1.93	0.84
1:A:186:GLU:OE1	1:A:187:LEU:HD23	1.78	0.83
1:C:24:LEU:HD12	1:C:50:TYR:HB2	1.59	0.83
1:D:132:ALA:HA	1:D:167:ILE:HD11	1.58	0.82
1:B:19:LEU:HD21	1:B:50:TYR:HB3	1.62	0.81
1:B:24:LEU:HD22	1:B:50:TYR:HB2	1.56	0.81
1:C:86:LEU:HB2	1:C:87:PRO:HA	1.62	0.80
1:B:97:ALA:CB	1:B:100:ARG:NH2	2.44	0.80
1:C:24:LEU:HD12	1:C:50:TYR:HB3	1.60	0.80
1:C:90:GLU:HG3	1:C:127:GLN:HG3	1.63	0.80
1:A:14:LEU:HD22	1:A:105:LEU:HD12	1.63	0.78
1:D:10:ARG:N	1:D:10:ARG:HD3	1.99	0.76
1:D:57:GLN:NE2	1:D:57:GLN:C	2.39	0.76
1:C:181:ARG:O	1:C:185:THR:HG23	1.84	0.76
1:D:9:LEU:HA	1:D:10:ARG:HH21	1.51	0.76
1:A:135:LEU:HD22	1:C:185:THR:HG21	1.66	0.76
1:B:69:GLN:C	1:B:71:HIS:O	2.25	0.75
1:A:97:ALA:HB3	1:A:100:ARG:HH11	1.51	0.75
2:A:201:COA:N6A	1:C:185:THR:HG22	2.02	0.75
1:B:43:GLY:O	1:B:49:TRP:CE3	2.41	0.74
1:A:135:LEU:CD2	1:C:185:THR:HG21	2.17	0.74

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:ASP:O	1:B:71:HIS:HB2	1.87	0.73
1:C:119:GLU:OE1	1:C:176:TRP:NE1	2.18	0.73
1:B:181:ARG:HH22	1:D:135:LEU:HB3	1.54	0.72
1:C:95:TRP:HA	2:C:201:COA:CDP	2.19	0.72
1:A:185:THR:HG1	1:A:186:GLU:N	1.88	0.71
1:D:137:ALA:O	1:D:141:ILE:HG13	1.90	0.71
1:A:28:LEU:HD22	1:A:47:PRO:HG3	1.73	0.71
1:A:14:LEU:HD22	1:A:105:LEU:CD1	2.19	0.71
1:B:43:GLY:H	1:B:44:PRO:HD3	1.56	0.70
1:A:176:TRP:HB3	1:A:177:PRO:HD3	1.73	0.70
1:D:24:LEU:HD21	1:D:47:PRO:HA	1.74	0.70
1:D:141:ILE:HD11	2:D:201:COA:H21	1.73	0.69
1:C:136:ARG:HH21	2:C:201:COA:H52A	1.58	0.69
1:C:138:GLN:O	1:C:142:GLU:HG3	1.93	0.69
1:B:92:GLY:HA2	2:B:201:COA:H32	1.74	0.69
1:B:97:ALA:CB	1:B:100:ARG:HH22	2.05	0.69
1:D:143:LYS:HE2	2:D:201:COA:O2A	1.92	0.68
1:A:16:LEU:HD21	1:A:109:ILE:HG23	1.75	0.68
1:D:101:HIS:NE2	2:D:201:COA:OAP	2.27	0.68
1:B:100:ARG:O	1:B:103:SER:OG	2.12	0.68
1:D:171:ILE:HD12	1:D:179:ILE:HG13	1.76	0.68
1:C:11:ARG:HH21	1:C:187:LEU:CD2	2.08	0.67
1:D:157:ARG:NH1	1:D:161:GLY:O	2.24	0.67
1:B:43:GLY:H	1:B:44:PRO:CD	2.07	0.66
1:A:82:PHE:CD1	1:A:89:VAL:HG23	2.31	0.66
1:C:96:LEU:HD21	1:C:109:ILE:HD11	1.77	0.66
1:B:59:GLU:N	1:B:59:GLU:OE1	2.28	0.66
1:A:151:GLN:NE2	1:A:165:ASP:OD2	2.28	0.65
1:C:86:LEU:CB	1:C:87:PRO:HA	2.21	0.65
1:D:148:ARG:HG3	1:D:169:TYR:CE2	2.31	0.65
1:B:68:ARG:HG3	1:B:73:LEU:CD2	2.24	0.65
1:D:14:LEU:HD13	1:D:105:LEU:HD21	1.79	0.65
1:B:38:LEU:CD2	1:B:95:TRP:HB2	2.27	0.65
1:B:24:LEU:HD21	1:B:50:TYR:HB2	0.65	0.64
1:B:183:LEU:O	1:B:187:LEU:HG	1.97	0.64
1:A:82:PHE:CE1	1:A:89:VAL:HG23	2.32	0.64
1:C:87:PRO:HD2	1:C:123:MET:HA	1.79	0.64
1:C:11:ARG:HE	1:C:187:LEU:HD21	1.62	0.64
1:A:178:GLU:OE1	1:A:181:ARG:HD2	1.97	0.64
1:D:43:GLY:O	1:D:45:LEU:N	2.31	0.63
1:B:174:ARG:NH2	1:D:154:ASN:OD1	2.31	0.63

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:130:THR:HB	2:D:201:COA:H31	1.81	0.63
1:B:19:LEU:HD13	1:B:64:ILE:HG12	1.81	0.62
1:A:180:ARG:O	1:A:184:GLU:HG3	1.99	0.62
1:B:14:LEU:HD11	1:B:67:VAL:HG13	1.82	0.62
1:B:89:VAL:HG13	1:B:123:MET:SD	2.40	0.62
1:A:133:SER:OG	1:A:164:ASP:OD2	2.14	0.62
1:D:19:LEU:HG	1:D:54:LEU:HD21	1.81	0.62
1:C:152:LEU:HD12	1:C:155:HIS:CE1	2.34	0.61
1:B:28:LEU:HD11	1:B:45:LEU:O	2.00	0.61
1:C:139:ARG:NH1	1:C:143:LYS:CG	2.64	0.61
1:D:130:THR:O	1:D:167:ILE:HD12	2.00	0.61
1:C:110:LYS:HD2	1:C:144:LEU:HD11	1.81	0.61
1:A:24:LEU:CD1	1:A:51:ARG:HG3	2.29	0.61
1:C:41:MET:O	1:C:159:ALA:CB	2.46	0.61
1:C:110:LYS:HB3	1:C:110:LYS:HZ2	1.61	0.60
1:B:181:ARG:HH12	1:D:135:LEU:HB2	1.67	0.60
2:B:201:COA:H131	2:B:201:COA:H72	1.84	0.60
1:D:46:ARG:HD2	1:D:47:PRO:HD2	1.84	0.60
1:D:143:LYS:CE	2:D:201:COA:O2A	2.49	0.60
1:B:43:GLY:O	1:B:49:TRP:CD2	2.55	0.59
1:D:135:LEU:HA	1:D:138:GLN:HB2	1.85	0.59
1:B:157:ARG:NH1	1:A:86:LEU:O	2.28	0.59
1:C:136:ARG:NH2	2:C:201:COA:H52A	2.17	0.58
1:C:64:ILE:HD12	1:C:65:TYR:H	1.68	0.58
2:B:201:COA:O9A	2:B:201:COA:O2B	2.18	0.58
1:A:185:THR:OG1	1:A:186:GLU:N	2.33	0.58
1:C:24:LEU:HD11	1:C:50:TYR:HB3	1.85	0.58
1:C:110:LYS:HB3	1:C:110:LYS:HZ3	1.66	0.58
1:B:86:LEU:HB2	1:B:87:PRO:HA	1.85	0.58
1:B:100:ARG:HH21	1:B:100:ARG:HG2	1.69	0.58
1:D:58:ARG:HB3	1:D:59:GLU:OE1	2.03	0.58
1:B:138:GLN:O	1:B:142:GLU:CG	2.49	0.58
1:A:130:THR:OG1	1:A:131:ALA:N	2.35	0.58
1:B:182:THR:O	1:B:185:THR:HB	2.04	0.57
1:A:157:ARG:NH1	1:A:161:GLY:O	2.35	0.57
1:B:26:GLU:OE2	1:B:71:HIS:ND1	2.37	0.57
1:B:28:LEU:HD13	1:B:47:PRO:HB3	1.87	0.57
1:B:183:LEU:C	1:B:185:THR:H	2.06	0.57
1:A:179:ILE:N	1:A:179:ILE:HD12	2.19	0.57
1:A:109:ILE:O	1:A:109:ILE:HG22	2.04	0.57
1:B:28:LEU:HD21	1:B:47:PRO:HD3	1.86	0.57

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:38:LEU:HD13	1:C:44:PRO:HB2	1.87	0.57
1:A:108:ALA:O	1:A:112:LEU:HG	2.04	0.57
1:A:175:ASP:C	1:A:179:ILE:HD13	2.17	0.56
1:C:110:LYS:CD	1:C:144:LEU:HD11	2.35	0.56
1:A:131:ALA:O	1:A:138:GLN:NE2	2.37	0.56
1:C:136:ARG:HE	2:C:201:COA:H51A	1.71	0.56
1:B:135:LEU:HD23	1:B:138:GLN:OE1	2.06	0.56
1:A:19:LEU:HB3	1:A:54:LEU:HD21	1.87	0.56
1:A:183:LEU:O	1:A:187:LEU:HD21	2.05	0.56
1:A:183:LEU:O	1:A:187:LEU:HG	2.06	0.56
1:D:57:GLN:C	1:D:57:GLN:HE21	2.07	0.56
1:A:24:LEU:HD11	1:A:54:LEU:HD12	1.88	0.56
1:A:178:GLU:O	1:A:182:THR:HG23	2.06	0.56
2:C:201:COA:H4B	2:C:201:COA:O2A	2.05	0.55
1:D:19:LEU:HB3	1:D:64:ILE:HG23	1.88	0.55
1:D:57:GLN:NE2	1:D:57:GLN:O	2.39	0.55
1:B:24:LEU:HD22	1:B:50:TYR:CD1	2.38	0.55
1:B:124:VAL:HA	1:B:173:ASP:HB3	1.88	0.55
1:C:84:PRO:C	1:C:86:LEU:H	2.09	0.55
1:B:28:LEU:CD2	1:B:47:PRO:HD3	2.35	0.55
1:C:2:PHE:HD1	1:C:3:LYS:N	2.05	0.55
1:D:70:ASP:OD2	1:D:71:HIS:ND1	2.40	0.55
1:C:64:ILE:HD13	1:C:78:ARG:NE	2.21	0.55
1:D:145:GLY:HA3	1:D:179:ILE:HD12	1.88	0.55
1:B:181:ARG:NH2	1:D:135:LEU:HB3	2.20	0.55
1:B:28:LEU:HD21	1:B:45:LEU:O	2.06	0.55
1:A:89:VAL:HB	1:A:123:MET:SD	2.47	0.54
1:B:26:GLU:CD	1:B:71:HIS:CE1	2.81	0.54
1:A:117:ALA:O	1:A:123:MET:HB2	2.06	0.54
1:A:19:LEU:CD2	1:A:64:ILE:HG12	2.38	0.54
1:A:10:ARG:NH1	1:A:12:GLY:O	2.40	0.54
1:A:24:LEU:HD21	1:A:54:LEU:HD11	1.90	0.54
1:A:134:ASN:HD21	2:A:201:COA:H71	1.72	0.54
1:A:182:THR:C	1:A:184:GLU:H	2.11	0.54
1:C:11:ARG:HH21	1:C:187:LEU:HD23	1.73	0.54
1:C:30:LEU:O	1:C:34:ASN:ND2	2.38	0.54
1:B:19:LEU:HD13	1:B:64:ILE:CG1	2.38	0.54
1:C:183:LEU:C	1:C:185:THR:H	2.10	0.54
1:B:153:ARG:NH1	1:A:170:SER:O	2.41	0.54
1:A:14:LEU:CD2	1:A:105:LEU:CD1	2.86	0.54
1:A:183:LEU:O	1:A:187:LEU:CG	2.56	0.54

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:86:LEU:CB	1:C:87:PRO:CA	2.85	0.54
1:C:110:LYS:HD3	1:C:144:LEU:HD21	1.89	0.54
1:B:174:ARG:HD3	1:D:165:ASP:OD1	2.08	0.53
1:A:14:LEU:CD2	1:A:105:LEU:HD12	2.36	0.53
1:A:131:ALA:HB3	1:A:134:ASN:HB2	1.89	0.53
2:D:201:COA:H143	2:D:201:COA:O4A	2.08	0.53
1:B:11:ARG:HB3	1:B:112:LEU:HD11	1.90	0.53
1:C:5:GLN:HG2	1:C:6:PRO:O	2.09	0.53
1:C:139:ARG:HH11	1:C:143:LYS:HG2	1.73	0.53
1:D:184:GLU:HA	1:D:187:LEU:HD12	1.90	0.53
1:B:32:GLU:HB2	1:B:45:LEU:HD23	1.91	0.53
1:C:63:VAL:HG23	1:C:79:PHE:HB2	1.91	0.53
1:C:99:ASP:HB2	1:C:100:ARG:NH2	2.24	0.53
1:C:43:GLY:HA3	1:C:49:TRP:CD1	2.44	0.53
1:D:118:PHE:O	1:D:122:GLU:HA	2.09	0.53
1:A:83:MET:HE3	1:A:86:LEU:HD12	1.91	0.53
1:C:69:GLN:O	1:C:72:ALA:N	2.31	0.53
1:C:76:THR:O	1:C:94:THR:HA	2.09	0.53
1:A:136:ARG:HG2	2:A:201:COA:H122	1.91	0.52
1:D:9:LEU:HB2	1:D:16:LEU:HD12	1.91	0.52
1:B:86:LEU:N	1:B:86:LEU:CD1	2.73	0.52
1:A:65:TYR:HB2	1:A:77:THR:O	2.10	0.52
1:D:50:TYR:O	1:D:54:LEU:HG	2.09	0.52
1:D:11:ARG:HB2	1:D:112:LEU:HD11	1.90	0.52
1:B:49:TRP:HZ3	1:B:95:TRP:HE1	1.57	0.51
1:C:91:ILE:HG21	1:C:113:MET:HE3	1.92	0.51
1:D:96:LEU:HD21	1:D:109:ILE:HD11	1.91	0.51
1:A:176:TRP:HB3	1:A:177:PRO:CD	2.40	0.51
1:C:177:PRO:O	1:C:181:ARG:HB2	2.10	0.51
1:D:24:LEU:HD11	1:D:47:PRO:HB3	1.92	0.51
1:B:44:PRO:HD2	1:B:45:LEU:H	1.74	0.51
1:A:114:LEU:CD2	1:A:126:VAL:HG11	2.40	0.51
1:B:87:PRO:HD2	1:B:123:MET:HA	1.93	0.51
1:B:110:LYS:HG3	1:B:144:LEU:HD11	1.92	0.51
1:A:11:ARG:NH2	1:A:187:LEU:HB3	2.26	0.51
1:D:147:ARG:NE	1:D:148:ARG:O	2.37	0.51
1:D:32:GLU:OE1	1:D:35:ARG:NH1	2.44	0.51
1:C:149:GLU:OE1	1:D:152:LEU:HD22	2.11	0.50
1:B:181:ARG:CZ	1:D:135:LEU:H	2.23	0.50
1:C:24:LEU:CD1	1:C:50:TYR:CB	2.65	0.50
1:D:24:LEU:HD23	1:D:50:TYR:HB2	1.92	0.50

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:182:THR:O	1:D:185:THR:OG1	2.22	0.50
1:D:182:THR:O	1:D:186:GLU:HG2	2.11	0.50
1:C:179:ILE:O	1:C:182:THR:HG22	2.11	0.50
1:D:89:VAL:HB	1:D:123:MET:SD	2.52	0.50
1:B:86:LEU:HD13	1:B:86:LEU:H	1.76	0.49
1:A:79:PHE:CE1	1:A:121:TRP:HZ3	2.30	0.49
1:C:178:GLU:O	1:C:179:ILE:HD13	2.11	0.49
1:B:16:LEU:HD21	1:B:67:VAL:HG22	1.93	0.49
1:B:154:ASN:HD22	1:A:124:VAL:HG22	1.77	0.49
1:A:14:LEU:HD11	1:A:105:LEU:HD11	1.89	0.49
2:A:201:COA:H61A	1:C:185:THR:HG22	1.76	0.49
1:D:11:ARG:NH1	1:D:187:LEU:HD22	2.28	0.49
1:B:69:GLN:O	1:B:71:HIS:O	2.31	0.49
1:A:24:LEU:HD22	1:A:50:TYR:HB3	1.94	0.49
1:A:14:LEU:CD1	1:A:105:LEU:CD1	2.73	0.49
1:B:111:PHE:HB2	1:B:183:LEU:HD13	1.93	0.49
1:A:182:THR:C	1:A:185:THR:OG1	2.49	0.49
1:C:21:GLU:HG3	1:C:51:ARG:HH22	1.78	0.48
1:C:94:THR:O	2:C:201:COA:H72	2.13	0.48
1:B:41:MET:O	1:B:41:MET:HG3	2.13	0.48
1:C:110:LYS:HE3	2:C:201:COA:S1P	2.53	0.48
1:B:16:LEU:CD2	1:B:67:VAL:HG22	2.42	0.48
1:C:24:LEU:N	1:C:24:LEU:HD22	2.28	0.48
1:C:156:ARG:NE	1:C:166:SER:OG	2.46	0.48
1:D:94:THR:O	2:D:201:COA:H62	2.14	0.48
1:A:26:GLU:HB3	1:A:73:LEU:HD13	1.95	0.48
1:A:54:LEU:O	1:A:57:GLN:HB3	2.13	0.48
1:B:26:GLU:O	1:B:26:GLU:HG2	2.13	0.48
1:C:158:LEU:HD12	1:C:162:ARG:NE	2.29	0.48
1:D:83:MET:O	1:D:87:PRO:HA	2.14	0.48
1:C:11:ARG:NH2	1:C:187:LEU:CD2	2.75	0.47
1:B:100:ARG:HD3	1:B:100:ARG:N	2.27	0.47
1:B:181:ARG:NH1	1:D:135:LEU:HB2	2.27	0.47
1:D:10:ARG:HD3	1:D:10:ARG:H	1.73	0.47
1:B:119:GLU:OE1	1:B:180:ARG:NH2	2.46	0.47
1:B:135:LEU:HA	1:B:138:GLN:OE1	2.14	0.47
1:B:151:GLN:HA	1:B:166:SER:O	2.15	0.47
1:B:183:LEU:O	1:B:185:THR:N	2.47	0.47
1:B:186:GLU:O	1:B:186:GLU:HG2	2.14	0.47
1:A:80:PHE:CD1	1:A:90:GLU:HB3	2.49	0.47
1:A:118:PHE:HE2	1:A:173:ASP:HA	1.79	0.47

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:27:LEU:HG	1:D:73:LEU:HD21	1.96	0.47
1:A:16:LEU:HD23	1:A:67:VAL:HG22	1.96	0.47
1:B:19:LEU:HD13	1:B:64:ILE:CD1	2.44	0.47
1:A:19:LEU:HD23	1:A:64:ILE:CD1	2.44	0.47
1:A:124:VAL:HG23	1:A:173:ASP:OD2	2.14	0.47
1:C:124:VAL:HG23	1:D:154:ASN:HB3	1.97	0.47
1:B:176:TRP:CZ2	1:B:180:ARG:HB2	2.50	0.47
1:B:158:LEU:HD12	1:B:162:ARG:HH11	1.79	0.47
1:A:182:THR:HA	1:A:185:THR:CG2	2.44	0.47
1:A:135:LEU:HD23	1:C:185:THR:HG21	1.94	0.47
1:C:7:VAL:HG23	1:C:65:TYR:OH	2.14	0.47
1:C:50:TYR:HH	1:C:95:TRP:HZ2	1.62	0.47
1:C:95:TRP:HA	2:C:201:COA:H132	1.95	0.47
1:A:118:PHE:CE2	1:A:173:ASP:HA	2.50	0.46
1:B:183:LEU:C	1:B:185:THR:N	2.68	0.46
1:C:11:ARG:NE	1:C:187:LEU:HD21	2.29	0.46
1:C:10:ARG:HG2	1:C:11:ARG:N	2.28	0.46
1:B:139:ARG:HA	1:B:142:GLU:HG3	1.97	0.46
2:B:201:COA:H131	2:B:201:COA:C7P	2.45	0.46
1:D:56:GLU:O	1:D:62:ALA:N	2.48	0.46
1:B:19:LEU:HD13	1:B:64:ILE:HD11	1.98	0.46
1:A:183:LEU:O	1:A:187:LEU:CD2	2.63	0.46
2:C:201:COA:N8P	2:C:201:COA:H133	2.30	0.46
1:A:110:LYS:HG3	1:A:144:LEU:HD21	1.98	0.46
1:C:134:ASN:O	1:C:136:ARG:N	2.48	0.46
1:C:158:LEU:CD1	1:C:162:ARG:NH1	2.78	0.46
1:A:98:GLU:HA	1:A:101:HIS:NE2	2.30	0.46
1:A:182:THR:OG1	1:A:183:LEU:N	2.49	0.46
1:C:70:ASP:O	1:C:71:HIS:ND1	2.48	0.46
1:A:82:PHE:CE1	1:A:89:VAL:CG2	2.99	0.46
1:B:152:LEU:HD22	1:A:168:LEU:HD13	1.98	0.45
1:C:158:LEU:HD23	1:C:158:LEU:HA	1.78	0.45
1:A:181:ARG:HD3	1:A:182:THR:CG2	2.47	0.45
1:D:31:ALA:HB1	1:D:45:LEU:HD23	1.98	0.45
1:C:109:ILE:HG22	1:C:113:MET:HE1	1.98	0.45
1:B:36:GLN:O	1:B:39:ILE:HG23	2.15	0.45
1:C:153:ARG:NH1	1:D:170:SER:O	2.49	0.45
1:B:162:ARG:HE	1:B:162:ARG:HB3	1.53	0.45
1:B:43:GLY:N	1:B:44:PRO:HD3	2.28	0.45
1:A:19:LEU:HG	1:A:54:LEU:HD21	1.99	0.45
1:C:57:GLN:HA	1:C:62:ALA:O	2.15	0.45

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:14:LEU:HD11	1:D:67:VAL:CG1	2.47	0.45
1:C:14:LEU:HD12	1:C:15:CYS:H	1.82	0.45
1:B:68:ARG:HG2	1:B:71:HIS:HA	1.97	0.45
1:B:176:TRP:CH2	1:B:180:ARG:HB2	2.52	0.45
1:C:130:THR:OG1	1:C:131:ALA:N	2.49	0.45
1:C:143:LYS:HD2	1:C:143:LYS:HA	1.70	0.44
1:B:71:HIS:HB3	1:B:72:ALA:HA	1.99	0.44
1:A:143:LYS:HE2	2:A:201:COA:H4B	1.98	0.44
1:A:36:GLN:O	1:A:39:ILE:HG12	2.18	0.44
1:C:156:ARG:HD2	1:C:165:ASP:O	2.16	0.44
1:B:28:LEU:HD11	1:B:45:LEU:C	2.38	0.44
1:B:170:SER:OG	1:A:153:ARG:NH1	2.44	0.44
1:C:28:LEU:O	1:C:32:GLU:N	2.43	0.44
1:C:42:ASN:HA	1:C:159:ALA:HB2	1.98	0.44
1:D:57:GLN:HE21	1:D:58:ARG:N	2.15	0.44
1:D:173:ASP:N	1:D:173:ASP:OD1	2.51	0.44
1:C:133:SER:O	1:C:135:LEU:N	2.48	0.44
1:D:100:ARG:HA	1:D:100:ARG:HD3	1.54	0.44
1:A:115:ARG:O	1:A:115:ARG:HG2	2.18	0.44
1:C:137:ALA:O	1:C:141:ILE:HG13	2.18	0.44
1:D:143:LYS:NZ	1:D:143:LYS:HB2	2.32	0.44
1:D:157:ARG:HB3	1:D:161:GLY:HA2	2.00	0.44
1:B:3:LYS:HE3	1:B:5:GLN:HE22	1.83	0.43
1:B:86:LEU:CB	1:B:87:PRO:CA	2.96	0.43
1:A:175:ASP:O	1:A:179:ILE:HD12	2.06	0.43
2:D:201:COA:H8A	2:D:201:COA:H2B	1.74	0.43
1:A:182:THR:HA	1:A:185:THR:HG23	1.99	0.43
1:A:182:THR:HA	1:A:185:THR:OG1	2.18	0.43
1:C:38:LEU:HB3	1:C:44:PRO:HG3	1.99	0.43
1:C:65:TYR:HB2	1:C:77:THR:O	2.19	0.43
1:B:86:LEU:HB2	1:B:87:PRO:CA	2.49	0.43
1:C:84:PRO:C	1:C:86:LEU:N	2.72	0.43
1:D:41:MET:SD	2:D:201:COA:H61	2.59	0.43
1:D:46:ARG:HH11	1:D:47:PRO:HD2	1.83	0.43
1:B:16:LEU:HD11	1:B:109:ILE:HG23	1.99	0.43
2:C:201:COA:O2A	2:C:201:COA:C4B	2.67	0.43
1:A:88:ALA:HA	1:A:123:MET:HG2	2.01	0.43
1:A:128:LEU:N	1:A:128:LEU:HD12	2.34	0.43
1:C:63:VAL:CG2	1:C:79:PHE:HB2	2.48	0.43
1:B:11:ARG:NH2	1:B:187:LEU:HB3	2.33	0.43
1:C:116:GLN:HA	1:C:120:GLN:HB2	1.99	0.43

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:183:LEU:O	1:C:185:THR:N	2.52	0.43
1:D:10:ARG:O	1:D:10:ARG:HG2	2.19	0.43
1:D:34:ASN:O	1:D:38:LEU:HG	2.18	0.43
1:A:53:GLY:HA3	1:A:64:ILE:CD1	2.49	0.43
1:C:152:LEU:HD12	1:C:155:HIS:NE2	2.33	0.43
1:A:182:THR:C	1:A:184:GLU:N	2.72	0.43
1:C:19:LEU:HB3	1:C:64:ILE:HG23	2.00	0.43
1:D:110:LYS:HA	1:D:110:LYS:HD3	1.78	0.43
1:C:124:VAL:HA	1:C:173:ASP:HB3	2.00	0.43
1:B:158:LEU:C	1:B:160:ASP:H	2.21	0.42
1:D:121:TRP:C	1:D:122:GLU:HG2	2.38	0.42
1:B:144:LEU:HD23	1:B:144:LEU:HA	1.77	0.42
1:B:154:ASN:ND2	1:A:124:VAL:HG22	2.34	0.42
1:A:55:ALA:HA	1:A:58:ARG:NH2	2.35	0.42
1:A:90:GLU:HG3	1:A:127:GLN:O	2.18	0.42
1:C:43:GLY:N	1:C:44:PRO:CD	2.83	0.42
1:C:109:ILE:CG2	1:C:113:MET:HE1	2.50	0.42
1:A:181:ARG:HD3	1:A:182:THR:HG22	2.01	0.42
1:A:111:PHE:CZ	1:A:184:GLU:OE2	2.72	0.42
1:A:118:PHE:HE1	1:A:126:VAL:HB	1.83	0.42
1:C:183:LEU:C	1:C:185:THR:N	2.73	0.42
1:D:144:LEU:HD23	1:D:144:LEU:HA	1.84	0.42
1:A:83:MET:O	1:A:87:PRO:HA	2.19	0.42
1:B:90:GLU:HG2	1:B:91:ILE:O	2.19	0.42
1:C:97:ALA:HB3	1:C:100:ARG:NH2	2.34	0.42
1:D:10:ARG:N	1:D:10:ARG:CD	2.71	0.42
1:B:3:LYS:HE3	1:B:5:GLN:NE2	2.35	0.42
1:A:48:ASP:HA	1:A:51:ARG:HB2	2.00	0.42
1:A:77:THR:HG21	1:A:109:ILE:HG21	2.01	0.42
1:C:3:LYS:HD2	1:C:3:LYS:HA	1.75	0.42
1:D:9:LEU:HA	1:D:10:ARG:NH2	2.28	0.42
2:A:201:COA:N3A	2:A:201:COA:H2B	2.34	0.42
1:D:3:LYS:HA	1:D:4:PRO:HD3	1.78	0.42
1:B:123:MET:HB2	1:B:123:MET:HE2	1.78	0.41
1:B:153:ARG:H	1:B:153:ARG:HG3	1.57	0.41
1:B:73:LEU:HD23	1:B:73:LEU:HA	1.81	0.41
1:B:38:LEU:HD21	1:B:95:TRP:CE3	2.55	0.41
1:B:82:PHE:CD1	1:B:89:VAL:HG12	2.54	0.41
1:A:117:ALA:O	1:A:121:TRP:HB2	2.21	0.41
1:C:98:GLU:C	1:C:100:ARG:H	2.24	0.41
1:C:101:HIS:CE1	2:C:201:COA:H143	2.56	0.41

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:158:LEU:HD12	1:C:162:ARG:HE	1.84	0.41
1:C:185:THR:H	1:C:185:THR:HG1	1.65	0.41
1:B:28:LEU:O	1:B:31:ALA:N	2.53	0.41
1:A:101:HIS:C	1:A:103:SER:H	2.24	0.41
2:A:201:COA:O7A	2:A:201:COA:C4B	2.68	0.41
1:D:185:THR:HG1	1:D:186:GLU:H	1.69	0.41
1:B:3:LYS:HE3	1:B:5:GLN:OE1	2.21	0.41
1:A:7:VAL:HG11	1:A:116:GLN:NE2	2.36	0.41
1:A:19:LEU:HD23	1:A:64:ILE:HD11	2.03	0.41
1:C:8:THR:HG22	1:C:18:PRO:HD3	2.03	0.41
1:D:89:VAL:HG22	1:D:90:GLU:N	2.36	0.41
1:D:101:HIS:CE1	2:D:201:COA:OAP	2.73	0.41
1:B:55:ALA:O	1:B:57:GLN:N	2.54	0.41
1:B:76:THR:O	1:B:94:THR:HA	2.21	0.41
1:B:79:PHE:O	1:B:80:PHE:HB3	2.21	0.41
1:B:147:ARG:O	1:B:169:TYR:HA	2.20	0.41
1:C:18:PRO:HA	1:C:64:ILE:O	2.21	0.41
1:C:64:ILE:HD13	1:C:78:ARG:HE	1.83	0.41
1:C:73:LEU:HA	1:C:73:LEU:HD13	1.83	0.41
1:C:109:ILE:HG22	1:C:113:MET:CE	2.51	0.41
1:C:141:ILE:O	1:C:144:LEU:HB2	2.21	0.41
1:A:105:LEU:O	1:A:109:ILE:HG13	2.21	0.40
1:C:135:LEU:HD23	1:C:135:LEU:HA	1.80	0.40
1:A:19:LEU:CB	1:A:54:LEU:HD21	2.50	0.40
1:A:79:PHE:CE1	1:A:121:TRP:CZ3	3.08	0.40
1:D:11:ARG:CZ	1:D:187:LEU:HD22	2.52	0.40
1:D:133:SER:OG	1:D:164:ASP:OD2	2.22	0.40
1:B:87:PRO:HB2	1:B:123:MET:HG2	2.03	0.40
1:A:53:GLY:HA3	1:A:64:ILE:HD13	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:ARG:NH2	1:A:188:SER:O[2_545]	2.19	0.01

## 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	185/193 (96%)	164 (89%)	18 (10%)	3 (2%)	9	37
1	B	185/193 (96%)	160 (86%)	17 (9%)	8 (4%)	2	14
1	C	185/193 (96%)	166 (90%)	14 (8%)	5 (3%)	5	24
1	D	185/193 (96%)	166 (90%)	14 (8%)	5 (3%)	5	24
All	All	740/772 (96%)	656 (89%)	63 (8%)	21 (3%)	5	23

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	44	PRO
1	B	56	GLU
1	C	86	LEU
1	D	4	PRO
1	B	71	HIS
1	C	184	GLU
1	B	4	PRO
1	B	55	ALA
1	A	70	ASP
1	C	134	ASN
1	C	135	LEU
1	D	31	ALA
1	D	44	PRO
1	D	70	ASP
1	B	184	GLU
1	A	86	LEU
1	A	98	GLU
1	B	86	LEU
1	D	86	LEU
1	C	74	VAL
1	B	43	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	154/157 (98%)	138 (90%)	16 (10%)	7	25
1	B	154/157 (98%)	142 (92%)	12 (8%)	12	39
1	C	154/157 (98%)	143 (93%)	11 (7%)	14	44
1	D	154/157 (98%)	146 (95%)	8 (5%)	23	57
All	All	616/628 (98%)	569 (92%)	47 (8%)	13	41

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	19	LEU
1	B	21	GLU
1	B	42	ASN
1	B	45	LEU
1	B	48	ASP
1	B	71	HIS
1	B	80	PHE
1	B	86	LEU
1	B	100	ARG
1	B	142	GLU
1	B	143	LYS
1	B	175	ASP
1	A	2	PHE
1	A	15	CYS
1	A	16	LEU
1	A	18	PRO
1	A	35	ARG
1	A	46	ARG
1	A	69	GLN
1	A	80	PHE
1	A	100	ARG
1	A	103	SER
1	A	143	LYS
1	A	153	ARG

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	170	SER
1	A	177	PRO
1	A	181	ARG
1	A	187	LEU
1	C	2	PHE
1	C	42	ASN
1	C	100	ARG
1	C	110	LYS
1	C	139	ARG
1	C	143	LYS
1	C	144	LEU
1	C	153	ARG
1	C	156	ARG
1	C	160	ASP
1	C	187	LEU
1	D	10	ARG
1	D	15	CYS
1	D	35	ARG
1	D	57	GLN
1	D	121	TRP
1	D	143	LYS
1	D	147	ARG
1	D	174	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	42	ASN
1	B	71	HIS
1	A	134	ASN
1	C	42	ASN
1	D	57	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	COA	B	201	-	41,50,50	4.19	16 (39%)	52,75,75	3.13	15 (28%)
2	COA	A	201	-	41,50,50	1.55	8 (19%)	52,75,75	2.57	17 (32%)
2	COA	D	201	-	41,50,50	4.25	19 (46%)	52,75,75	3.05	15 (28%)
2	COA	C	201	-	41,50,50	2.02	15 (36%)	52,75,75	2.65	15 (28%)

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	COA	B	201	-	-	24/44/64/64	0/3/3/3
2	COA	A	201	-	-	26/44/64/64	0/3/3/3
2	COA	D	201	-	-	22/44/64/64	0/3/3/3
2	COA	C	201	-	-	16/44/64/64	0/3/3/3

All (58) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	201	COA	O4B-C1B	16.36	1.63	1.41
2	B	201	COA	O4B-C1B	16.10	1.63	1.41
2	B	201	COA	C2B-C1B	-13.92	1.32	1.53
2	D	201	COA	C2B-C1B	-13.45	1.33	1.53
2	D	201	COA	C9P-N8P	10.06	1.55	1.33
2	B	201	COA	C9P-N8P	9.64	1.54	1.33

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	201	COA	O4B-C4B	-6.37	1.30	1.45
2	B	201	COA	O4B-C4B	-6.36	1.30	1.45
2	D	201	COA	C6P-C5P	5.21	1.61	1.51
2	A	201	COA	C4A-N3A	-4.45	1.29	1.35
2	D	201	COA	O5P-C5P	-4.30	1.14	1.23
2	B	201	COA	O5P-C5P	-4.27	1.14	1.23
2	C	201	COA	O5P-C5P	-3.87	1.15	1.23
2	B	201	COA	C6P-C5P	3.84	1.58	1.51
2	B	201	COA	C5P-N4P	3.73	1.41	1.33
2	D	201	COA	C7P-N8P	3.70	1.54	1.46
2	C	201	COA	O4B-C1B	-3.59	1.36	1.41
2	C	201	COA	P1A-O1A	-3.47	1.38	1.50
2	A	201	COA	CEP-CBP	-3.29	1.46	1.53
2	C	201	COA	C3B-C4B	-3.27	1.44	1.52
2	B	201	COA	C7P-N8P	3.17	1.53	1.46
2	C	201	COA	C2A-N3A	-3.15	1.27	1.32
2	C	201	COA	C4A-N3A	-3.10	1.31	1.35
2	D	201	COA	C5P-N4P	3.07	1.40	1.33
2	C	201	COA	P1A-O2A	-2.99	1.41	1.55
2	C	201	COA	O9P-C9P	-2.90	1.17	1.23
2	B	201	COA	O2B-C2B	2.83	1.49	1.43
2	D	201	COA	O3B-C3B	-2.76	1.34	1.44
2	A	201	COA	P2A-O5A	-2.66	1.42	1.55
2	D	201	COA	OAP-CAP	-2.66	1.37	1.42
2	D	201	COA	C5A-C4A	-2.65	1.33	1.40
2	A	201	COA	P3B-O9A	-2.65	1.44	1.54
2	B	201	COA	CEP-CBP	2.65	1.59	1.53
2	B	201	COA	O3B-C3B	-2.60	1.34	1.44
2	B	201	COA	C5A-C4A	-2.55	1.34	1.40
2	A	201	COA	P3B-O8A	-2.53	1.45	1.54
2	B	201	COA	C6A-N6A	2.52	1.43	1.34
2	C	201	COA	C2A-N1A	-2.52	1.29	1.33
2	B	201	COA	C3B-C4B	2.50	1.59	1.52
2	D	201	COA	CEP-CBP	2.47	1.59	1.53
2	D	201	COA	C6A-N6A	2.45	1.43	1.34
2	C	201	COA	P1A-O5B	-2.45	1.49	1.59
2	C	201	COA	CEP-CBP	-2.41	1.48	1.53
2	B	201	COA	CDP-CBP	2.38	1.59	1.53
2	A	201	COA	C5A-C4A	2.37	1.47	1.40
2	D	201	COA	O2B-C2B	2.29	1.48	1.43
2	A	201	COA	C5P-N4P	-2.27	1.28	1.33
2	D	201	COA	C7P-C6P	2.22	1.58	1.51

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	201	COA	P3B-O3B	2.21	1.63	1.59
2	D	201	COA	C3B-C4B	2.20	1.58	1.52
2	D	201	COA	CDP-CBP	2.19	1.58	1.53
2	C	201	COA	P2A-O5A	-2.19	1.45	1.55
2	C	201	COA	P3B-O9A	-2.18	1.46	1.54
2	C	201	COA	C5A-N7A	-2.18	1.31	1.39
2	A	201	COA	C5A-N7A	-2.14	1.32	1.39
2	D	201	COA	P2A-O6A	2.13	1.67	1.59
2	C	201	COA	O4B-C4B	-2.13	1.40	1.45
2	B	201	COA	P3B-O3B	2.13	1.63	1.59

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	201	COA	C5A-C6A-N6A	13.34	140.62	120.35
2	D	201	COA	C5A-C6A-N6A	13.21	140.42	120.35
2	A	201	COA	CDP-CBP-CAP	9.40	125.13	108.82
2	B	201	COA	N6A-C6A-N1A	-9.28	99.31	118.57
2	D	201	COA	N6A-C6A-N1A	-9.18	99.51	118.57
2	C	201	COA	O4B-C1B-C2B	-7.80	95.53	106.93
2	B	201	COA	CAP-C9P-N8P	7.36	131.24	116.58
2	A	201	COA	CEP-CBP-CAP	-7.29	96.17	108.82
2	C	201	COA	CEP-CBP-CAP	-6.63	97.32	108.82
2	D	201	COA	C7P-C6P-C5P	6.55	123.26	112.36
2	C	201	COA	C2B-C3B-C4B	-6.40	91.89	103.22
2	D	201	COA	O6A-CCP-CBP	6.35	120.76	110.55
2	C	201	COA	N3A-C2A-N1A	-5.99	119.32	128.68
2	D	201	COA	N3A-C2A-N1A	-5.89	119.47	128.68
2	B	201	COA	N3A-C2A-N1A	-5.87	119.51	128.68
2	A	201	COA	P2A-O3A-P1A	-5.03	115.56	132.83
2	C	201	COA	O3B-C3B-C4B	-4.85	92.54	110.08
2	B	201	COA	O6A-CCP-CBP	-4.78	102.87	110.55
2	B	201	COA	O9P-C9P-N8P	-4.71	112.87	122.99
2	C	201	COA	CDP-CBP-CCP	4.56	115.67	108.23
2	B	201	COA	CDP-CBP-CAP	4.31	116.30	108.82
2	A	201	COA	C3P-N4P-C5P	-4.27	114.91	122.84
2	C	201	COA	C6P-C5P-N4P	4.26	123.59	116.42
2	B	201	COA	P2A-O3A-P1A	-4.04	118.95	132.83
2	D	201	COA	C7P-N8P-C9P	3.97	129.68	122.59
2	C	201	COA	CAP-C9P-N8P	3.95	124.44	116.58
2	B	201	COA	C7P-N8P-C9P	3.83	129.42	122.59
2	A	201	COA	O6A-CCP-CBP	3.77	116.60	110.55

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	201	COA	O5P-C5P-C6P	-3.75	115.16	122.02
2	A	201	COA	C7P-N8P-C9P	3.69	129.17	122.59
2	D	201	COA	O5P-C5P-N4P	-3.67	116.10	123.01
2	C	201	COA	C1B-N9A-C4A	-3.55	120.40	126.64
2	A	201	COA	CEP-CBP-CDP	-3.52	101.99	109.17
2	C	201	COA	C2A-N1A-C6A	3.46	124.67	118.75
2	A	201	COA	CAP-C9P-N8P	3.41	123.36	116.58
2	A	201	COA	CEP-CBP-CCP	3.29	113.60	108.23
2	D	201	COA	CDP-CBP-CAP	3.28	114.50	108.82
2	A	201	COA	C7P-C6P-C5P	3.20	117.68	112.36
2	B	201	COA	CDP-CBP-CCP	-3.03	103.30	108.23
2	A	201	COA	CDP-CBP-CCP	-2.99	103.35	108.23
2	A	201	COA	O9A-P3B-O7A	2.80	121.63	110.68
2	C	201	COA	C3P-N4P-C5P	-2.74	117.75	122.84
2	D	201	COA	CEP-CBP-CAP	-2.71	104.12	108.82
2	C	201	COA	C3B-C2B-C1B	2.67	105.80	99.89
2	B	201	COA	O4B-C1B-C2B	-2.65	103.05	106.93
2	D	201	COA	OAP-CAP-CBP	-2.59	104.16	110.25
2	D	201	COA	CAP-C9P-N8P	2.58	121.71	116.58
2	D	201	COA	C6P-C5P-N4P	2.51	120.64	116.42
2	D	201	COA	CEP-CBP-CCP	2.50	112.31	108.23
2	C	201	COA	O9A-P3B-O8A	2.46	117.04	107.64
2	A	201	COA	N3A-C2A-N1A	-2.44	124.87	128.68
2	C	201	COA	O9P-C9P-CAP	-2.38	113.79	121.06
2	B	201	COA	OAP-CAP-CBP	-2.38	104.65	110.25
2	B	201	COA	C6P-C5P-N4P	2.38	120.42	116.42
2	B	201	COA	CEP-CBP-CCP	-2.36	104.38	108.23
2	A	201	COA	C6P-C7P-N8P	-2.29	107.27	111.90
2	A	201	COA	N6A-C6A-N1A	2.28	123.32	118.57
2	B	201	COA	O5P-C5P-N4P	-2.20	118.86	123.01
2	A	201	COA	C3B-C2B-C1B	2.19	104.73	99.89
2	D	201	COA	C1B-N9A-C4A	2.14	130.41	126.64
2	D	201	COA	P2A-O3A-P1A	-2.03	125.85	132.83
2	A	201	COA	P2A-O6A-CCP	2.01	133.17	121.56

There are no chirality outliers.

All (88) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	201	COA	C5B-O5B-P1A-O3A
2	B	201	COA	CCP-O6A-P2A-O3A
2	B	201	COA	CCP-O6A-P2A-O5A

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms
2	B	201	COA	CAP-CBP-CCP-O6A
2	B	201	COA	OAP-CAP-CBP-CCP
2	B	201	COA	C9P-CAP-CBP-CCP
2	B	201	COA	OAP-CAP-CBP-CDP
2	B	201	COA	C9P-CAP-CBP-CDP
2	B	201	COA	OAP-CAP-CBP-CEP
2	B	201	COA	C9P-CAP-CBP-CEP
2	B	201	COA	O9P-C9P-CAP-CBP
2	B	201	COA	N8P-C9P-CAP-CBP
2	B	201	COA	N8P-C9P-CAP-OAP
2	B	201	COA	CAP-C9P-N8P-C7P
2	B	201	COA	S1P-C2P-C3P-N4P
2	A	201	COA	C4B-C3B-O3B-P3B
2	A	201	COA	C5B-O5B-P1A-O3A
2	A	201	COA	CAP-CBP-CCP-O6A
2	A	201	COA	OAP-CAP-CBP-CCP
2	A	201	COA	C9P-CAP-CBP-CCP
2	A	201	COA	OAP-CAP-CBP-CDP
2	A	201	COA	C9P-CAP-CBP-CDP
2	A	201	COA	OAP-CAP-CBP-CEP
2	A	201	COA	C9P-CAP-CBP-CEP
2	A	201	COA	O9P-C9P-CAP-CBP
2	A	201	COA	N8P-C9P-CAP-CBP
2	A	201	COA	N8P-C9P-CAP-OAP
2	A	201	COA	C6P-C5P-N4P-C3P
2	A	201	COA	O5P-C5P-N4P-C3P
2	A	201	COA	S1P-C2P-C3P-N4P
2	C	201	COA	C3B-O3B-P3B-O7A
2	C	201	COA	C4B-C5B-O5B-P1A
2	C	201	COA	CEP-CBP-CCP-O6A
2	C	201	COA	CAP-CBP-CCP-O6A
2	C	201	COA	C9P-CAP-CBP-CCP
2	C	201	COA	C9P-CAP-CBP-CEP
2	C	201	COA	O9P-C9P-CAP-CBP
2	C	201	COA	N8P-C9P-CAP-CBP
2	C	201	COA	C6P-C5P-N4P-C3P
2	C	201	COA	O5P-C5P-N4P-C3P
2	C	201	COA	S1P-C2P-C3P-N4P
2	D	201	COA	C5B-O5B-P1A-O3A
2	D	201	COA	CBP-CCP-O6A-P2A
2	D	201	COA	CDP-CBP-CCP-O6A
2	D	201	COA	CEP-CBP-CCP-O6A

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms
2	D	201	COA	CAP-CBP-CCP-O6A
2	D	201	COA	OAP-CAP-CBP-CCP
2	D	201	COA	C9P-CAP-CBP-CCP
2	D	201	COA	OAP-CAP-CBP-CDP
2	D	201	COA	C9P-CAP-CBP-CDP
2	D	201	COA	OAP-CAP-CBP-CEP
2	D	201	COA	C9P-CAP-CBP-CEP
2	D	201	COA	O9P-C9P-CAP-CBP
2	D	201	COA	N8P-C9P-CAP-CBP
2	D	201	COA	N8P-C9P-CAP-OAP
2	D	201	COA	C5P-C6P-C7P-N8P
2	D	201	COA	S1P-C2P-C3P-N4P
2	B	201	COA	O9P-C9P-N8P-C7P
2	B	201	COA	O4B-C4B-C5B-O5B
2	A	201	COA	C3B-C4B-C5B-O5B
2	A	201	COA	O4B-C4B-C5B-O5B
2	C	201	COA	CDP-CBP-CCP-O6A
2	A	201	COA	O9P-C9P-CAP-OAP
2	C	201	COA	O9P-C9P-CAP-OAP
2	B	201	COA	CEP-CBP-CCP-O6A
2	A	201	COA	CAP-C9P-N8P-C7P
2	A	201	COA	O9P-C9P-N8P-C7P
2	D	201	COA	O4B-C4B-C5B-O5B
2	D	201	COA	P2A-O3A-P1A-O1A
2	D	201	COA	P1A-O3A-P2A-O5A
2	B	201	COA	C5B-O5B-P1A-O1A
2	B	201	COA	C5B-O5B-P1A-O2A
2	B	201	COA	CCP-O6A-P2A-O4A
2	D	201	COA	C5B-O5B-P1A-O2A
2	B	201	COA	C3B-C4B-C5B-O5B
2	D	201	COA	C2P-C3P-N4P-C5P
2	C	201	COA	OAP-CAP-CBP-CCP
2	B	201	COA	CDP-CBP-CCP-O6A
2	A	201	COA	CEP-CBP-CCP-O6A
2	A	201	COA	C5P-C6P-C7P-N8P
2	A	201	COA	C2P-C3P-N4P-C5P
2	B	201	COA	O5P-C5P-C6P-C7P
2	C	201	COA	C3B-O3B-P3B-O9A
2	C	201	COA	O4B-C4B-C5B-O5B
2	D	201	COA	P1A-O3A-P2A-O4A
2	A	201	COA	C5B-O5B-P1A-O1A
2	A	201	COA	CCP-O6A-P2A-O4A

*Continued on next page...*

*Continued from previous page...*

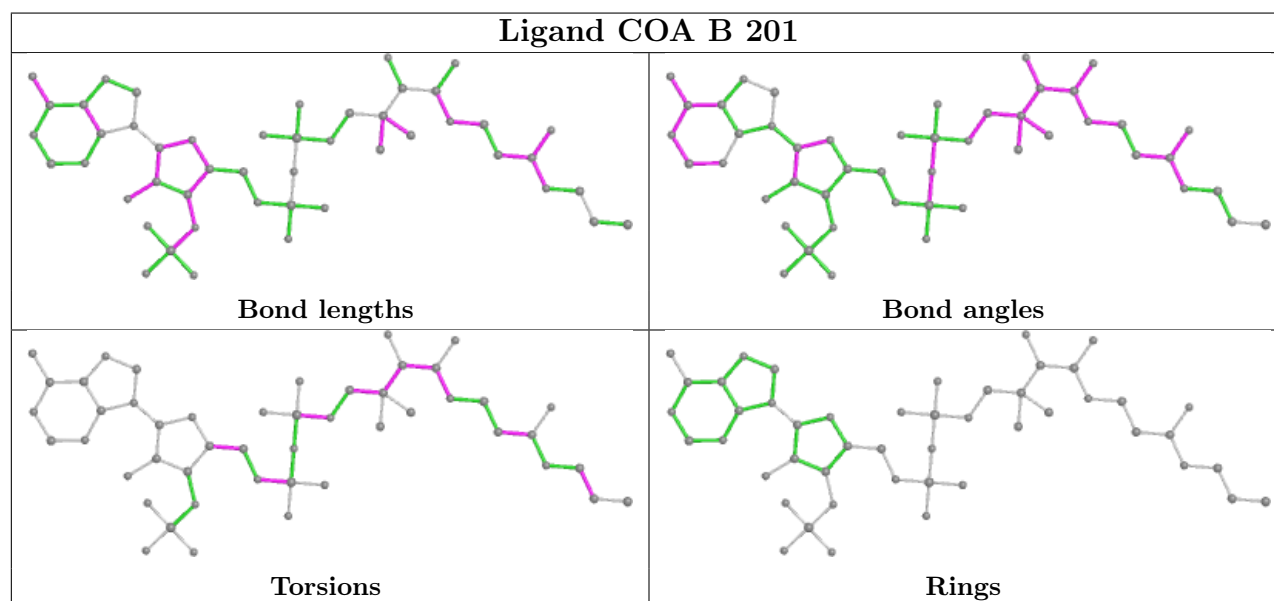
Mol	Chain	Res	Type	Atoms
2	A	201	COA	C2B-C3B-O3B-P3B

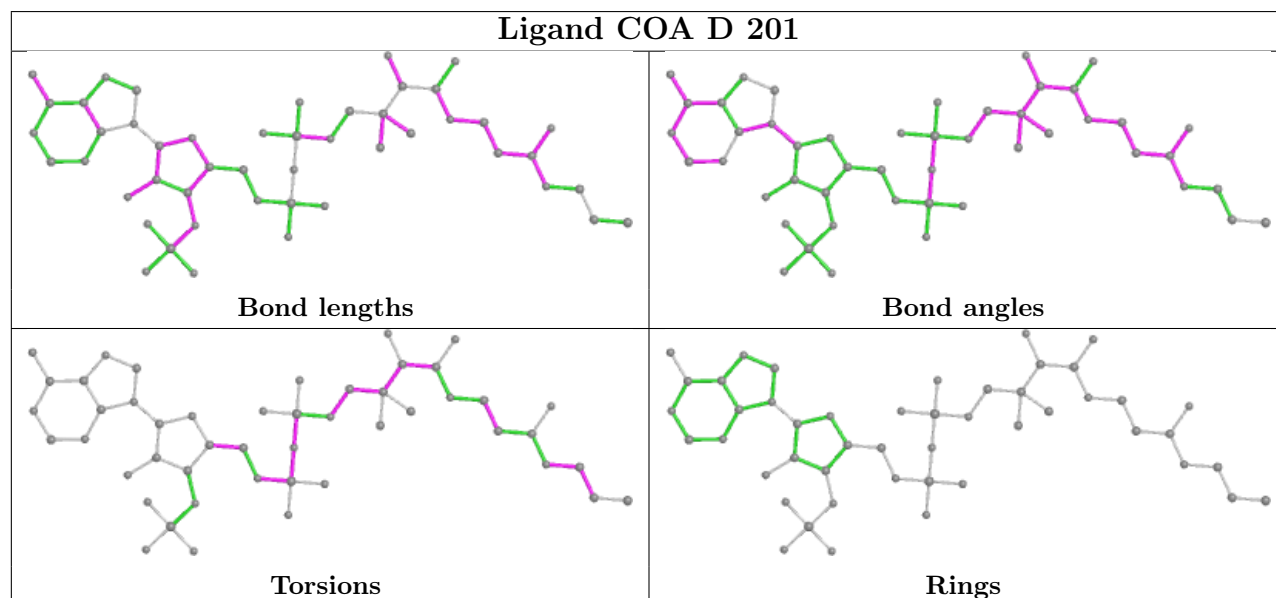
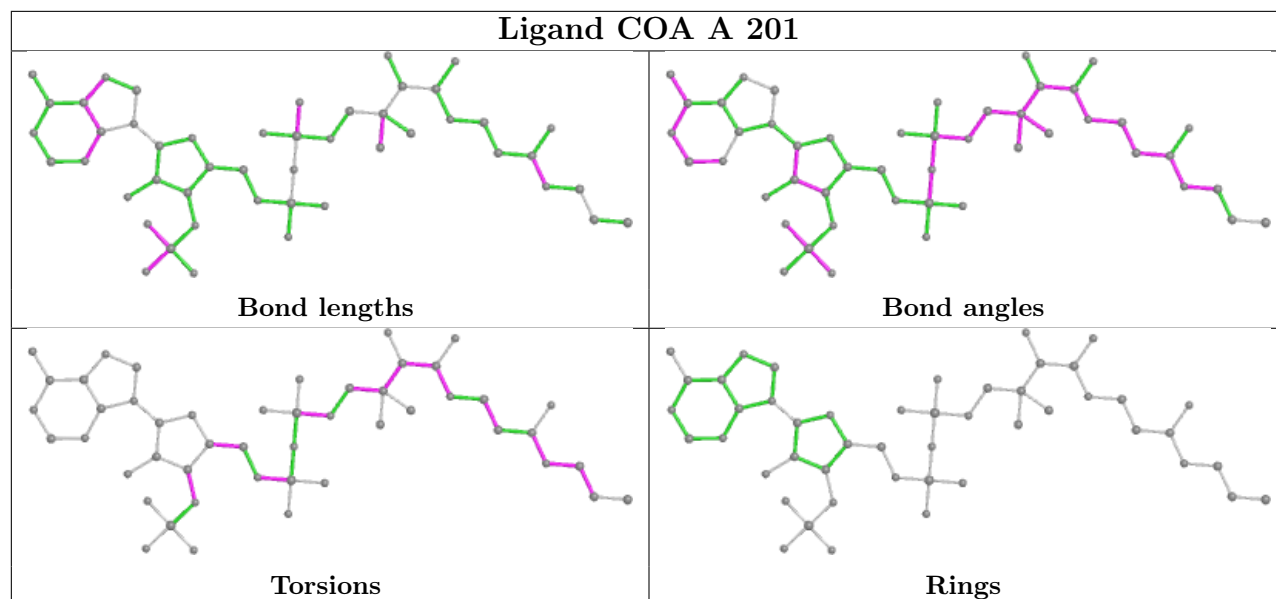
There are no ring outliers.

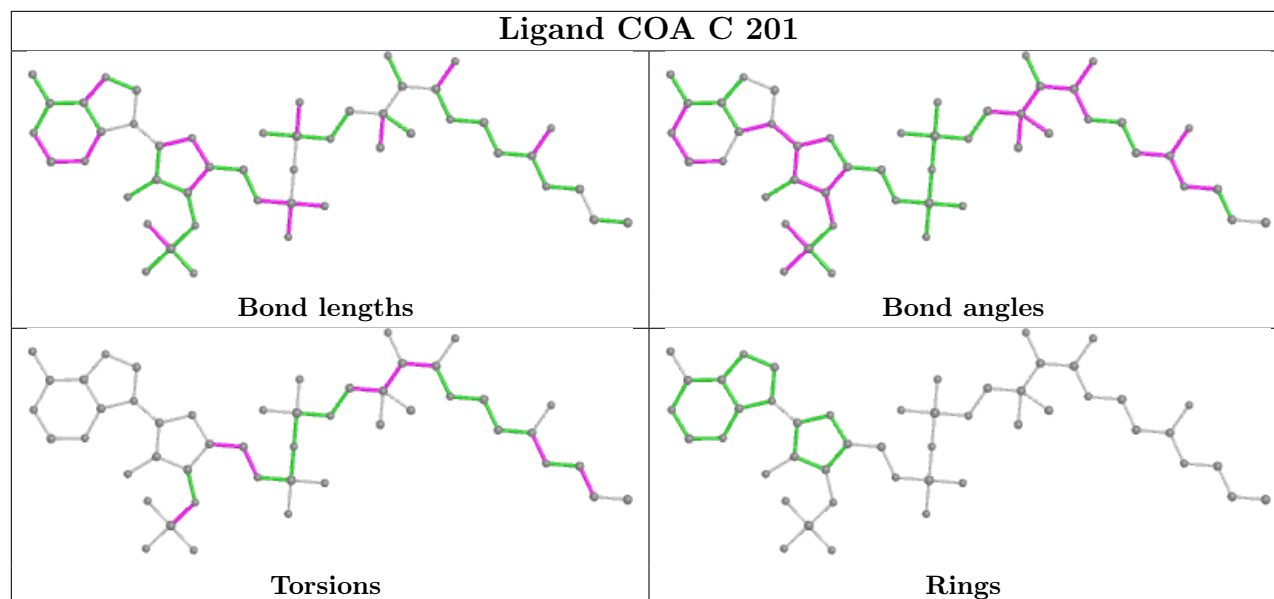
4 monomers are involved in 34 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	201	COA	5	0
2	A	201	COA	7	0
2	D	201	COA	11	0
2	C	201	COA	11	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 [i](#)

There are no chain breaks in this entry.



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

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	187/193 (96%)	-0.31	0 100 100	13, 26, 41, 55	0
1	B	187/193 (96%)	-0.20	0 100 100	14, 25, 48, 55	0
1	C	187/193 (96%)	-0.20	1 (0%) 91 75	12, 27, 43, 62	0
1	D	187/193 (96%)	-0.10	2 (1%) 80 56	13, 26, 46, 58	0
All	All	748/772 (96%)	-0.20	3 (0%) 92 79	12, 26, 47, 62	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	70	ASP	2.5
1	D	160	ASP	2.3
1	C	48	ASP	2.3

### 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](#)

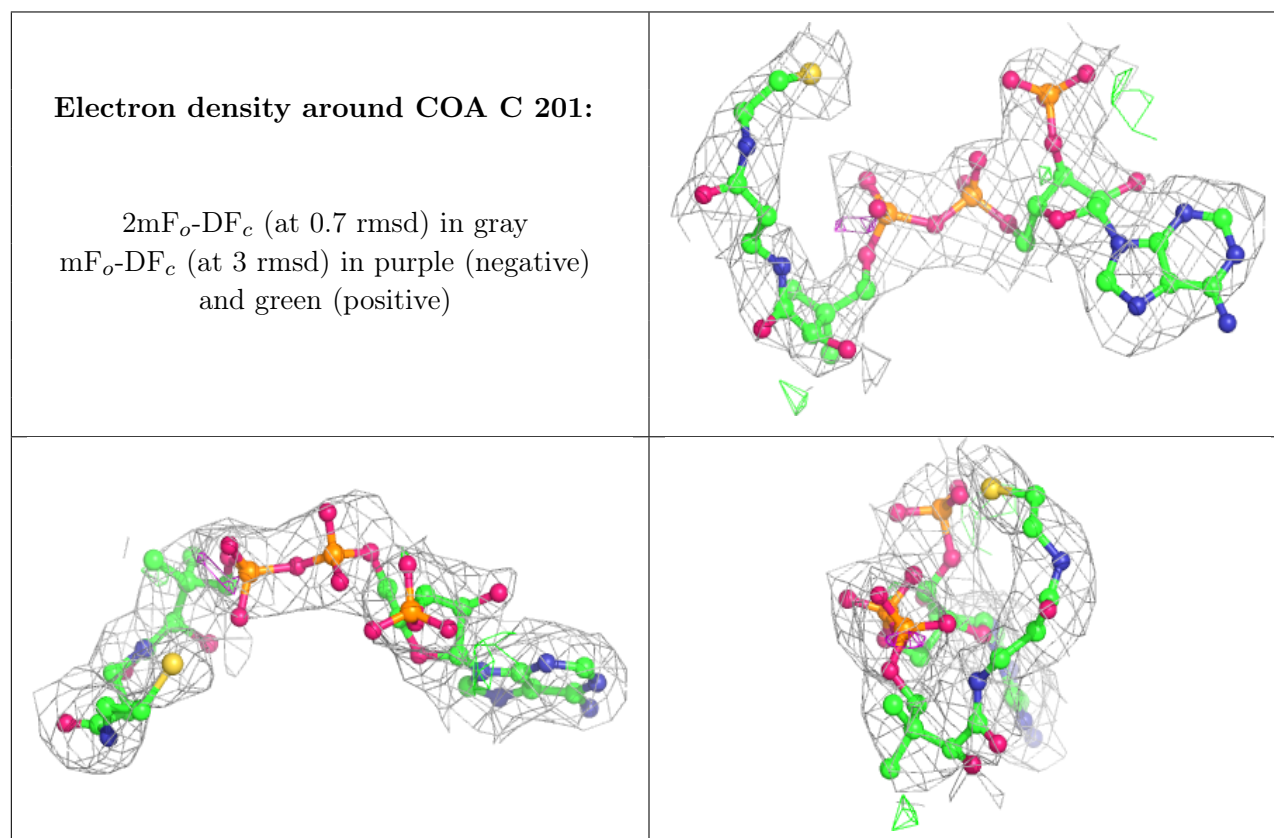
There are no monosaccharides in this entry.

### 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, 95<sup>th</sup> 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.

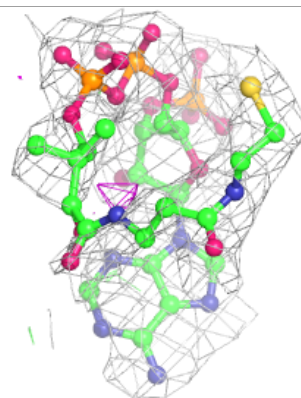
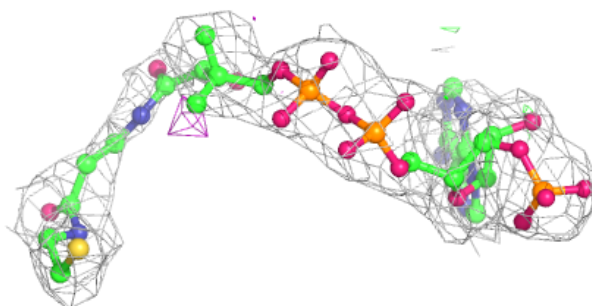
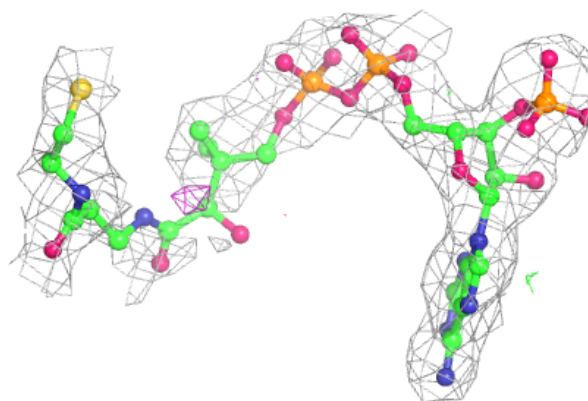
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	COA	C	201	48/48	0.86	0.23	22,40,52,63	0
2	COA	A	201	48/48	0.89	0.22	26,49,68,78	0
2	COA	B	201	48/48	0.91	0.19	14,29,44,49	0
2	COA	D	201	48/48	0.92	0.22	15,28,38,43	0

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.

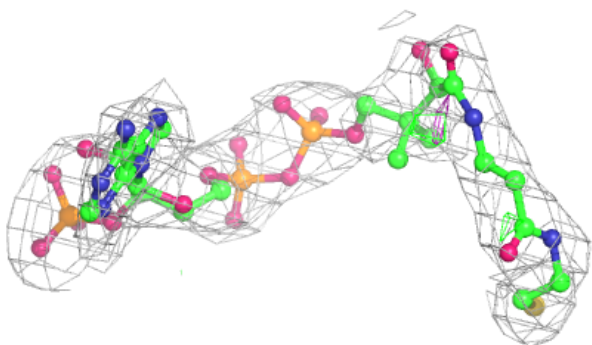
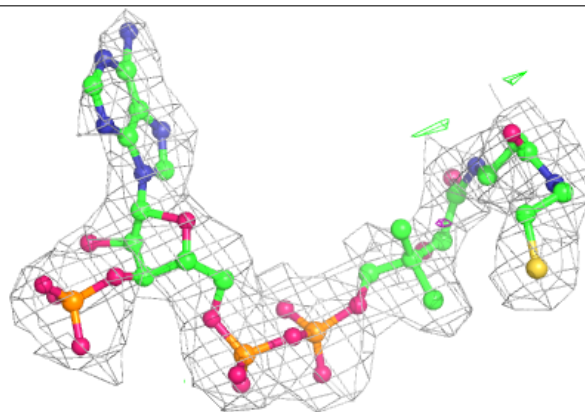


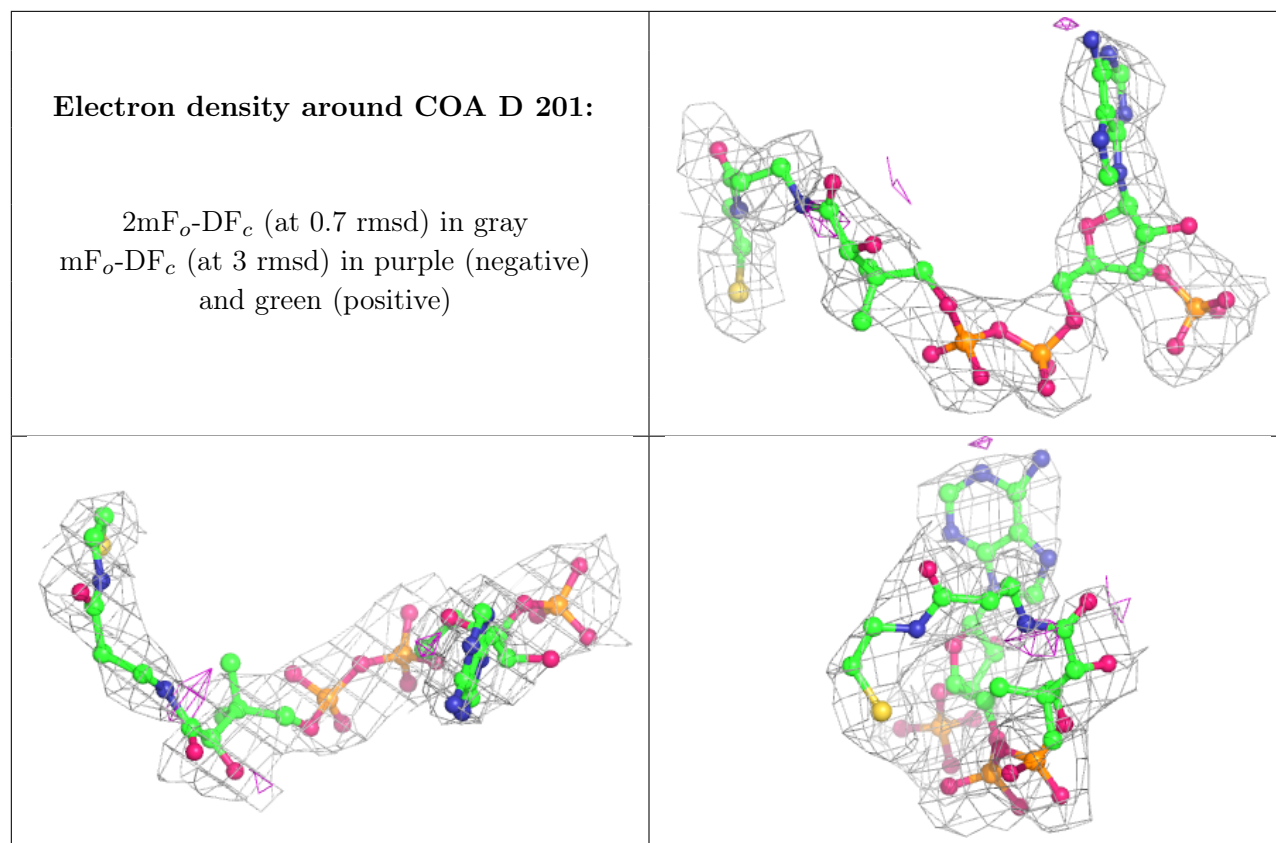
**Electron density around COA A 201:**

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

**Electron density around COA B 201:**

$2mF_o-DF_c$  (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.