



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:10 PM UTC

PDB ID : 2WPW / pdb_00002wpw
Title : Tandem GNAT protein from the clavulanic acid biosynthesis pathway (without AcCoA)
Authors : Iqbal, A.; Arunlanantham, H.; McDonough, M.A.; Chowdhury, R.; Clifton, I.J.
Deposited on : 2009-08-11
Resolution : 2.38 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)
Xtrriage (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

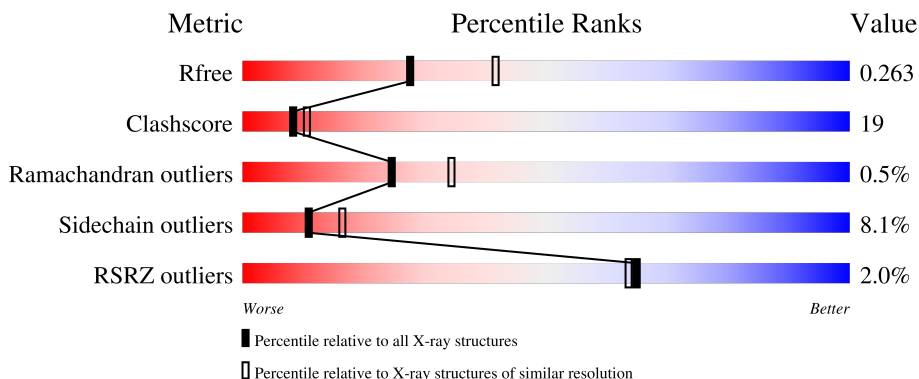
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.38 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	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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	339	 2% 71% 21% . . .
1	B	339	 2% 69% 22% . .
1	C	339	 % 73% 18% 5% .
1	D	339	 % 68% 24% . .

2 Entry composition [i](#)

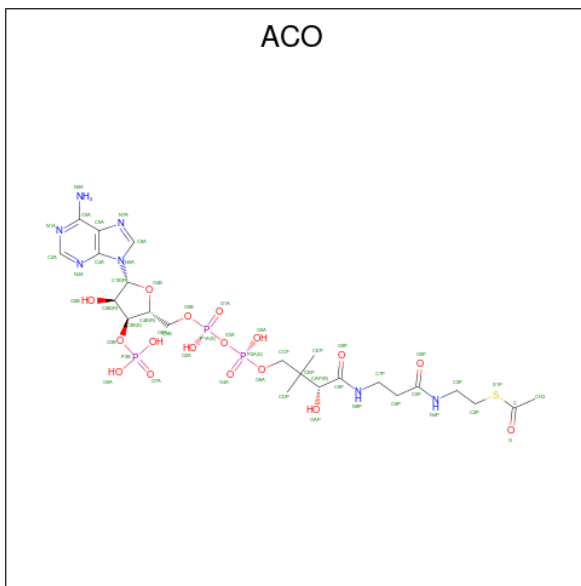
There are 3 unique types of molecules in this entry. The entry contains 11259 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ORF14.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	328	Total 2486	C 1554	N 462	O 463	S 1	Se 6	0	0	0
1	B	326	Total 2474	C 1547	N 460	O 460	S 1	Se 6	0	0	0
1	C	325	Total 2470	C 1545	N 459	O 459	S 1	Se 6	0	0	0
1	D	325	Total 2470	C 1545	N 459	O 459	S 1	Se 6	0	0	0

- Molecule 2 is ACETYL COENZYME *A (CCD ID: ACO) (formula: C₂₃H₃₈N₇O₁₇P₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	N	O	P	S		
2	A	1	Total 51	C 23	N 7	O 17	P 3	S 1	0	0
2	B	1	Total 51	C 23	N 7	O 17	P 3	S 1	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
2	C	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		
2	D	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		

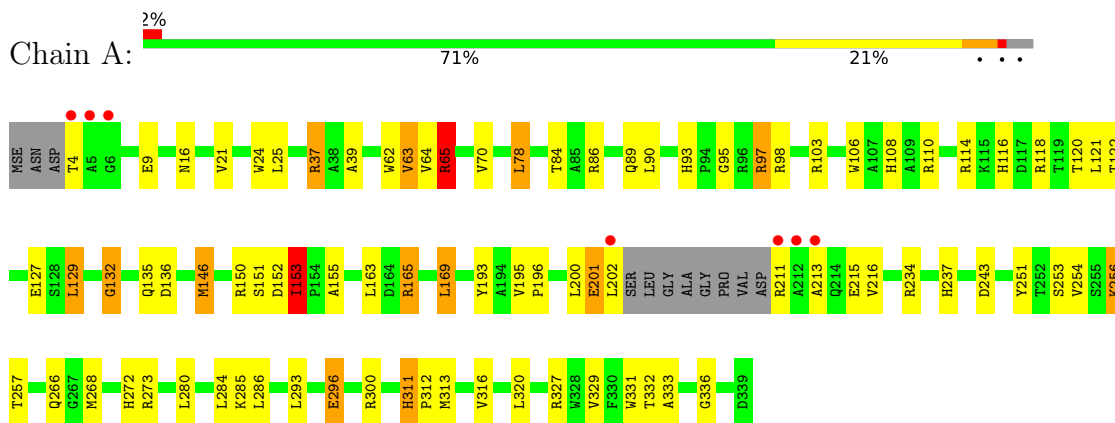
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	285	Total	O	0	0
			285	285		
3	B	277	Total	O	0	0
			277	277		
3	C	283	Total	O	0	0
			283	283		
3	D	310	Total	O	0	0
			310	310		

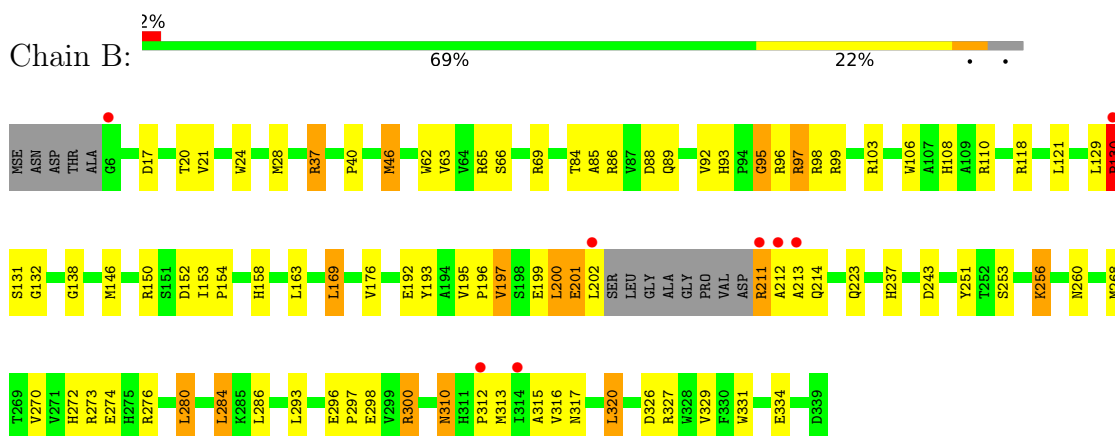
3 Residue-property plots [i](#)

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.

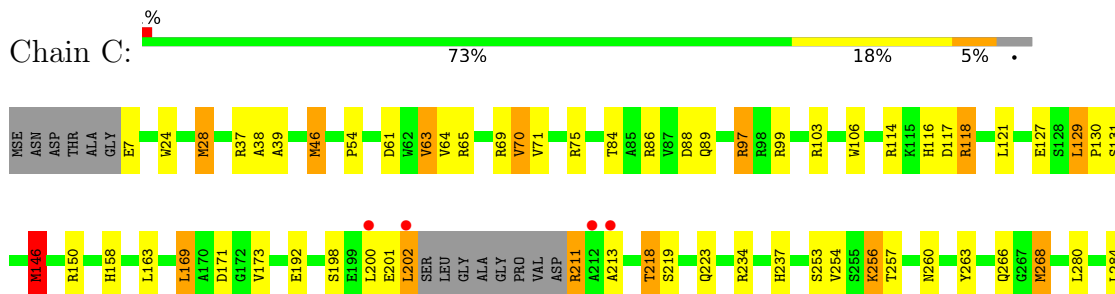
- Molecule 1: ORF14



- Molecule 1: ORF14

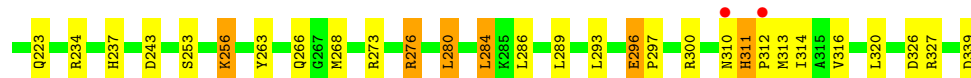
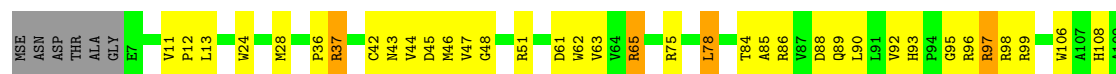


- Molecule 1: ORF14





• Molecule 1: ORF14



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.45Å 93.50Å 126.43Å 90.00° 97.86° 90.00°	Depositor
Resolution (Å)	32.40 – 2.38 32.40 – 2.38	Depositor EDS
% Data completeness (in resolution range)	92.2 (32.40-2.38) 92.4 (32.40-2.38)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 2.39Å)	Xtrriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.209 , 0.266 0.214 , 0.263	Depositor DCC
R_{free} test set	3318 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	30.0	Xtrriage
Anisotropy	0.287	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11259	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.35 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.3381e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: ACO

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	1.11	1/2543 (0.0%)	1.18	9/3468 (0.3%)
1	B	1.10	3/2531 (0.1%)	1.19	10/3451 (0.3%)
1	C	1.12	3/2527 (0.1%)	1.17	12/3446 (0.3%)
1	D	1.09	0/2527	1.15	9/3446 (0.3%)
All	All	1.10	7/10128 (0.1%)	1.17	40/13811 (0.3%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	28	MSE	SE-CE	-7.31	1.73	1.95
1	B	46	MSE	SE-CE	-6.20	1.76	1.95
1	A	132	GLY	N-CA	6.09	1.51	1.43
1	C	146	MSE	SE-CE	-5.67	1.78	1.95
1	C	173	VAL	N-CA	-5.60	1.41	1.46
1	B	153	ILE	CA-C	5.16	1.57	1.52
1	B	40	PRO	CA-C	5.11	1.54	1.51

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	39	ALA	CA-C-N	-8.36	111.77	120.38
1	C	39	ALA	C-N-CA	-8.36	111.77	120.38
1	B	260	ASN	CA-C-N	8.36	128.85	119.32
1	B	260	ASN	C-N-CA	8.36	128.85	119.32
1	B	153	ILE	CA-C-N	-8.24	111.30	119.78
1	B	153	ILE	C-N-CA	-8.24	111.30	119.78
1	A	132	GLY	CA-C-N	8.03	128.50	119.83
1	A	132	GLY	C-N-CA	8.03	128.50	119.83
1	B	138	GLY	CA-C-N	-7.95	110.22	119.05
1	B	138	GLY	C-N-CA	-7.95	110.22	119.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	132	GLY	CA-C-N	7.82	127.58	119.76
1	D	132	GLY	C-N-CA	7.82	127.58	119.76
1	C	296	GLU	CA-C-N	6.53	126.22	119.56
1	C	296	GLU	C-N-CA	6.53	126.22	119.56
1	D	176	VAL	CB-CA-C	-6.44	103.27	110.68
1	A	311	HIS	CA-C-N	6.21	125.43	118.97
1	A	311	HIS	C-N-CA	6.21	125.43	118.97
1	C	173	VAL	CA-C-N	5.96	126.30	119.92
1	C	173	VAL	C-N-CA	5.96	126.30	119.92
1	A	39	ALA	N-CA-C	-5.85	102.21	109.64
1	C	260	ASN	CA-C-N	5.84	126.25	119.47
1	C	260	ASN	C-N-CA	5.84	126.25	119.47
1	A	65	ARG	N-CA-C	5.72	118.73	109.40
1	B	197	VAL	CB-CA-C	-5.70	104.68	111.97
1	B	176	VAL	CB-CA-C	-5.67	105.67	110.94
1	D	311	HIS	CA-C-N	5.56	126.79	119.84
1	D	311	HIS	C-N-CA	5.56	126.79	119.84
1	A	153	ILE	CA-C-N	-5.55	114.21	119.76
1	A	153	ILE	C-N-CA	-5.55	114.21	119.76
1	B	95	GLY	N-CA-C	-5.48	106.24	115.62
1	D	136	ASP	CA-C-N	-5.48	114.37	120.45
1	D	136	ASP	C-N-CA	-5.48	114.37	120.45
1	D	197	VAL	CB-CA-C	-5.41	104.95	112.04
1	C	336	GLY	CA-C-N	5.31	125.78	120.52
1	C	336	GLY	C-N-CA	5.31	125.78	120.52
1	C	70	VAL	N-CA-C	5.24	116.00	108.93
1	A	155	ALA	N-CA-C	5.20	116.80	110.41
1	D	65	ARG	N-CA-C	5.05	117.05	109.23
1	B	243	ASP	N-CA-C	5.03	117.14	111.11
1	C	46	MSE	N-CA-C	5.01	116.55	111.14

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	2486	0	2420	92	0
1	B	2474	0	2408	92	0
1	C	2470	0	2405	79	0
1	D	2470	0	2405	102	0
2	A	51	0	34	5	0
2	B	51	0	34	5	0
2	C	51	0	34	4	1
2	D	51	0	34	8	1
3	A	285	0	0	30	0
3	B	277	0	0	21	0
3	C	283	0	0	31	0
3	D	310	0	0	50	0
All	All	11259	0	9774	382	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 19.

All (382) 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:103:ARG:HA	1:B:146:MSE:CE	1.59	1.33
1:D:313:MSE:HG2	3:D:2048:HOH:O	1.30	1.24
1:A:132:GLY:HA2	3:A:2127:HOH:O	1.39	1.21
1:D:150:ARG:NH1	3:D:2154:HOH:O	1.76	1.17
2:A:401:ACO:C8A	2:A:401:ACO:H51A	1.76	1.15
1:D:276:ARG:HG2	1:D:276:ARG:HH11	1.13	1.13
1:C:337:PRO:O	3:C:2273:HOH:O	1.65	1.13
1:B:103:ARG:HA	1:B:146:MSE:HE1	1.23	1.11
2:B:401:ACO:H52A	2:B:401:ACO:C8A	1.79	1.11
2:C:401:ACO:H51A	2:C:401:ACO:H8A	1.33	1.11
1:D:273:ARG:HG3	3:D:2263:HOH:O	1.50	1.10
2:B:401:ACO:H52A	2:B:401:ACO:H8A	1.13	1.09
2:C:401:ACO:H51A	2:C:401:ACO:C8A	1.84	1.07
1:D:44:VAL:N	3:D:2054:HOH:O	1.79	1.07
1:C:312:PRO:O	1:C:316:VAL:HG23	1.50	1.07
1:A:200:LEU:O	1:A:201:GLU:HB2	1.54	1.06
1:D:36:PRO:O	3:D:2034:HOH:O	1.70	1.06
1:D:201:GLU:O	1:D:202:LEU:HD12	1.55	1.06
2:D:401:ACO:H8A	2:D:401:ACO:H51A	1.13	1.06
1:A:256:LYS:HA	3:A:2213:HOH:O	1.55	1.05
1:A:257:THR:N	3:A:2213:HOH:O	1.81	1.04
1:B:103:ARG:HA	1:B:146:MSE:HE2	1.37	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:401:ACO:H51A	2:A:401:ACO:H8A	1.09	1.03
1:A:165:ARG:HB3	1:A:165:ARG:HH11	1.21	1.01
1:C:316:VAL:HG11	3:C:2236:HOH:O	1.57	1.01
1:C:296:GLU:HG3	3:C:2242:HOH:O	1.61	1.00
1:B:93:HIS:CD2	1:B:95:GLY:H	1.79	1.00
1:D:234:ARG:HD3	3:D:2234:HOH:O	1.58	1.00
1:D:243:ASP:OD2	3:D:2238:HOH:O	1.81	0.97
2:D:401:ACO:H51A	2:D:401:ACO:C8A	1.93	0.97
1:B:28:MSE:SE	1:B:46:MSE:HE2	2.15	0.96
1:B:103:ARG:CA	1:B:146:MSE:HE1	1.95	0.96
1:B:93:HIS:HD2	1:B:95:GLY:H	0.97	0.94
1:D:150:ARG:NH2	3:D:2154:HOH:O	2.00	0.94
1:C:106:TRP:HB3	1:C:146:MSE:HE2	1.46	0.94
1:D:316:VAL:HG11	3:D:2273:HOH:O	1.67	0.93
1:C:171:ASP:HB2	3:C:2156:HOH:O	1.67	0.93
2:D:401:ACO:H8A	2:D:401:ACO:C5B	1.99	0.92
1:C:65:ARG:HD2	3:C:2076:HOH:O	1.69	0.91
1:C:316:VAL:CG1	3:C:2236:HOH:O	2.15	0.91
1:A:336:GLY:HA3	3:A:2109:HOH:O	1.70	0.90
1:A:110:ARG:NH2	3:A:2109:HOH:O	2.05	0.89
1:D:93:HIS:CD2	1:D:95:GLY:H	1.91	0.89
1:A:273:ARG:NH1	1:A:273:ARG:HB2	1.88	0.88
1:D:200:LEU:O	1:D:201:GLU:HB2	1.74	0.87
2:B:401:ACO:H8A	2:B:401:ACO:C5B	2.03	0.87
1:A:201:GLU:HB3	3:A:2179:HOH:O	1.75	0.85
1:C:266:GLN:OE1	3:C:2221:HOH:O	1.93	0.85
1:D:28:MSE:SE	1:D:46:MSE:HE2	2.25	0.85
1:A:106:TRP:HB3	1:A:146:MSE:HE2	1.57	0.84
1:D:93:HIS:HD2	1:D:95:GLY:H	1.21	0.84
1:D:46:MSE:HE1	3:D:2088:HOH:O	1.76	0.84
1:A:243:ASP:HB2	3:A:2206:HOH:O	1.76	0.83
1:D:150:ARG:CZ	3:D:2154:HOH:O	2.10	0.83
1:B:93:HIS:HD2	1:B:95:GLY:N	1.77	0.82
2:D:401:ACO:O8A	3:D:2308:HOH:O	1.97	0.82
1:B:96:ARG:HD3	1:B:99:ARG:NH1	1.95	0.81
1:A:165:ARG:HH11	1:A:165:ARG:CB	1.92	0.81
1:B:66:SER:O	3:B:2072:HOH:O	1.96	0.81
1:A:211:ARG:HG3	1:A:213:ALA:H	1.44	0.81
1:A:24:TRP:HE1	1:A:89:GLN:HE22	1.28	0.81
1:D:276:ARG:HH11	1:D:276:ARG:CG	1.89	0.81
1:B:106:TRP:HB3	1:B:146:MSE:HE3	1.62	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:ARG:HD2	3:A:2067:HOH:O	1.80	0.80
1:A:122:THR:HG22	1:A:332:THR:HG22	1.64	0.80
1:C:24:TRP:HE1	1:C:89:GLN:HE22	1.27	0.80
1:D:24:TRP:HE1	1:D:89:GLN:HE22	1.27	0.79
1:B:24:TRP:HE1	1:B:89:GLN:HE22	1.30	0.79
1:B:103:ARG:HG3	1:B:146:MSE:HE1	1.63	0.78
1:D:276:ARG:HG2	1:D:276:ARG:NH1	1.95	0.78
1:A:65:ARG:NH2	3:A:2066:HOH:O	2.15	0.78
1:D:311:HIS:N	1:D:312:PRO:CD	2.47	0.78
1:D:96:ARG:HD2	1:D:99:ARG:NH1	1.99	0.77
1:A:256:LYS:CA	3:A:2213:HOH:O	2.18	0.77
1:B:212:ALA:C	1:B:214:GLN:H	1.90	0.77
1:B:192:GLU:HG2	3:B:2183:HOH:O	1.85	0.77
1:D:310:ASN:C	1:D:312:PRO:HD2	2.09	0.77
2:D:401:ACO:C8A	2:D:401:ACO:C5B	2.61	0.76
1:B:202:LEU:HD23	3:B:2232:HOH:O	1.84	0.76
1:A:93:HIS:HD2	1:A:95:GLY:H	1.32	0.76
1:D:37:ARG:HD2	3:D:2039:HOH:O	1.86	0.76
2:A:401:ACO:C8A	2:A:401:ACO:C5B	2.61	0.76
1:D:98:ARG:NH2	3:D:2109:HOH:O	2.18	0.76
1:B:97:ARG:HB3	2:B:401:ACO:O4A	1.85	0.75
1:B:331:TRP:O	3:B:2261:HOH:O	2.03	0.75
1:A:234:ARG:HD2	3:A:2201:HOH:O	1.85	0.75
1:D:193:TYR:C	1:D:196:PRO:HD2	2.12	0.75
1:A:78:LEU:HD21	1:A:116:HIS:CD2	2.22	0.74
1:B:297:PRO:HD2	3:B:2246:HOH:O	1.86	0.74
1:C:234:ARG:NH2	1:C:296:GLU:OE1	2.20	0.74
1:D:154:PRO:O	3:D:2160:HOH:O	2.04	0.74
1:C:127:GLU:OE2	1:C:150:ARG:NH1	2.20	0.73
1:D:276:ARG:NH1	3:D:2266:HOH:O	1.90	0.73
1:A:106:TRP:CD1	1:A:146:MSE:HE3	2.25	0.72
1:A:234:ARG:CD	3:A:2201:HOH:O	2.38	0.71
1:A:273:ARG:CB	1:A:273:ARG:HH11	2.03	0.71
1:C:103:ARG:HG3	1:C:146:MSE:HE1	1.71	0.71
1:D:218:THR:HG22	3:D:2216:HOH:O	1.89	0.71
1:D:88:ASP:HB3	3:D:2100:HOH:O	1.90	0.71
1:C:234:ARG:HD3	3:C:2200:HOH:O	1.91	0.71
1:D:266:GLN:HG2	3:D:2255:HOH:O	1.89	0.71
1:C:46:MSE:HE1	3:C:2081:HOH:O	1.92	0.70
1:C:211:ARG:HG2	3:C:2181:HOH:O	1.91	0.70
1:B:96:ARG:HD3	1:B:99:ARG:CZ	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:ARG:HD2	3:A:2029:HOH:O	1.93	0.68
1:B:46:MSE:HE1	3:B:2079:HOH:O	1.93	0.68
1:B:103:ARG:CG	1:B:146:MSE:HE1	2.22	0.68
2:C:401:ACO:H8A	2:C:401:ACO:C5B	2.19	0.68
1:D:93:HIS:HD2	1:D:95:GLY:N	1.91	0.68
2:D:401:ACO:N1A	3:D:2309:HOH:O	2.26	0.67
1:A:84:THR:O	1:A:118:ARG:HG2	1.93	0.67
1:B:103:ARG:CA	1:B:146:MSE:CE	2.53	0.67
1:C:313:MSE:CG	3:C:2115:HOH:O	2.43	0.67
1:C:268:MSE:HE3	3:C:2224:HOH:O	1.95	0.66
1:D:169:LEU:HD13	1:D:286:LEU:HD22	1.76	0.66
1:D:237:HIS:CD2	1:D:253:SER:HB3	2.31	0.66
1:B:97:ARG:O	1:B:98:ARG:HB2	1.95	0.66
1:D:78:LEU:HD21	1:D:116:HIS:CD2	2.30	0.66
1:B:202:LEU:HD21	1:B:276:ARG:NH2	2.11	0.66
1:C:86:ARG:HG3	3:C:2123:HOH:O	1.96	0.66
1:B:202:LEU:CD2	3:B:2232:HOH:O	2.43	0.65
1:A:146:MSE:CE	3:A:2105:HOH:O	2.43	0.65
1:A:106:TRP:CG	1:A:146:MSE:HG3	2.32	0.65
1:C:106:TRP:CG	1:C:146:MSE:HG3	2.32	0.65
1:A:313:MSE:HG2	3:A:2259:HOH:O	1.96	0.65
1:C:211:ARG:HD2	1:C:213:ALA:H	1.63	0.64
1:B:69:ARG:HB3	3:B:2072:HOH:O	1.97	0.64
1:B:212:ALA:C	1:B:214:GLN:N	2.53	0.64
1:C:106:TRP:CD1	1:C:146:MSE:HE3	2.33	0.63
1:A:234:ARG:NH2	1:A:296:GLU:OE1	2.30	0.63
1:B:272:HIS:CE1	1:B:274:GLU:H	2.15	0.62
2:A:401:ACO:H8A	2:A:401:ACO:C5B	2.05	0.62
1:B:106:TRP:HB3	1:B:146:MSE:CE	2.28	0.62
1:A:106:TRP:HD1	1:A:146:MSE:HE3	1.65	0.62
1:C:65:ARG:CD	3:C:2076:HOH:O	2.34	0.62
1:C:106:TRP:HB3	1:C:146:MSE:CE	2.26	0.61
1:D:296:GLU:HG3	3:D:2251:HOH:O	2.00	0.61
1:A:336:GLY:CA	3:A:2109:HOH:O	2.38	0.61
1:D:300:ARG:CG	3:D:2079:HOH:O	2.49	0.61
1:A:195:VAL:HG23	3:A:2177:HOH:O	2.00	0.61
1:A:273:ARG:HB2	1:A:273:ARG:HH11	1.55	0.61
1:B:300:ARG:NH2	1:B:300:ARG:HG3	2.16	0.61
1:A:243:ASP:CB	3:A:2206:HOH:O	2.40	0.61
1:D:234:ARG:NH2	1:D:296:GLU:OE1	2.27	0.61
1:B:310:ASN:ND2	1:B:310:ASN:O	2.33	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:211:ARG:CG	3:C:2181:HOH:O	2.48	0.61
1:C:63:VAL:HG11	1:C:70:VAL:HG22	1.82	0.60
1:D:90:LEU:C	1:D:90:LEU:HD23	2.26	0.60
1:D:110:ARG:NH1	3:D:2124:HOH:O	2.21	0.60
1:D:276:ARG:NH2	3:D:2266:HOH:O	2.34	0.60
1:B:129:LEU:O	1:B:131:SER:N	2.34	0.60
1:C:103:ARG:HA	1:C:146:MSE:CE	2.31	0.60
1:C:106:TRP:CB	1:C:146:MSE:HE2	2.25	0.60
1:C:313:MSE:HG2	3:C:2115:HOH:O	2.01	0.60
1:D:300:ARG:HG2	3:D:2079:HOH:O	2.01	0.60
1:D:276:ARG:CG	1:D:276:ARG:NH1	2.57	0.60
1:B:154:PRO:O	3:B:2146:HOH:O	2.17	0.60
1:C:256:LYS:HE2	1:C:263:TYR:O	2.01	0.60
1:C:311:HIS:O	1:C:312:PRO:C	2.45	0.60
1:A:97:ARG:HB3	2:A:401:ACO:O4A	2.02	0.59
1:B:298:GLU:O	1:B:300:ARG:HD2	2.02	0.59
1:C:38:ALA:HB2	3:C:2044:HOH:O	2.03	0.59
1:C:198:SER:HA	3:C:2178:HOH:O	2.03	0.59
1:A:63:VAL:HG11	1:A:70:VAL:HG22	1.84	0.59
1:B:195:VAL:HB	1:B:196:PRO:HD3	1.85	0.59
1:A:146:MSE:HE3	3:A:2105:HOH:O	2.02	0.59
1:A:84:THR:C	1:A:118:ARG:HG2	2.28	0.58
1:D:129:LEU:HD13	1:D:327:ARG:HB2	1.85	0.58
1:B:199:GLU:C	1:B:201:GLU:N	2.60	0.58
1:C:54:PRO:O	3:C:2062:HOH:O	2.16	0.58
1:B:256:LYS:HD2	1:B:296:GLU:OE1	2.04	0.58
1:A:165:ARG:CB	1:A:165:ARG:NH1	2.66	0.58
1:B:199:GLU:C	1:B:201:GLU:H	2.10	0.58
1:D:51:ARG:HD3	3:D:2058:HOH:O	2.03	0.58
1:B:280:LEU:O	1:B:284:LEU:HD22	2.04	0.57
1:C:103:ARG:CG	1:C:146:MSE:HE1	2.34	0.57
1:A:254:VAL:HG23	3:A:2241:HOH:O	2.04	0.57
1:C:84:THR:O	1:C:118:ARG:HG2	2.05	0.57
1:A:84:THR:HG22	1:A:120:THR:OG1	2.04	0.57
1:C:256:LYS:HD3	1:C:256:LYS:C	2.30	0.57
1:D:97:ARG:O	1:D:98:ARG:HB2	2.03	0.57
1:D:311:HIS:N	1:D:312:PRO:HD2	2.16	0.57
1:B:193:TYR:O	1:B:197:VAL:HG23	2.05	0.57
1:A:62:TRP:CD2	1:A:108:HIS:HE1	2.23	0.57
1:D:280:LEU:O	1:D:284:LEU:HD22	2.05	0.57
1:C:201:GLU:O	1:C:202:LEU:HD12	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:268:MSE:HE1	1:D:313:MSE:HE1	1.87	0.56
1:D:268:MSE:HE3	3:D:2261:HOH:O	2.04	0.56
1:A:273:ARG:NH1	1:A:273:ARG:CB	2.61	0.56
1:A:311:HIS:N	1:A:312:PRO:CD	2.67	0.56
1:B:202:LEU:HB2	1:B:270:VAL:CG1	2.35	0.56
1:C:28:MSE:HE1	1:C:46:MSE:SE	2.55	0.56
1:A:127:GLU:OE2	1:A:150:ARG:NH1	2.37	0.56
1:A:169:LEU:CD1	1:A:286:LEU:HD22	2.36	0.56
1:B:86:ARG:HD2	1:B:88:ASP:OD2	2.05	0.56
1:D:339:ASP:CG	3:D:2304:HOH:O	2.48	0.56
1:C:114:ARG:HD2	1:C:336:GLY:O	2.06	0.56
1:C:310:ASN:CG	1:C:313:MSE:SE	2.98	0.55
1:C:313:MSE:HG3	3:C:2115:HOH:O	2.04	0.55
1:D:219:SER:O	1:D:223:GLN:HG3	2.07	0.55
1:C:169:LEU:HD22	1:C:289:LEU:HD23	1.89	0.55
1:A:193:TYR:C	1:A:196:PRO:HD2	2.32	0.55
1:C:256:LYS:HA	3:C:2209:HOH:O	2.06	0.55
1:A:327:ARG:HD3	3:A:2268:HOH:O	2.06	0.55
1:D:313:MSE:HB3	3:D:2117:HOH:O	2.07	0.55
1:B:315:ALA:HA	3:B:2255:HOH:O	2.05	0.54
1:D:84:THR:O	1:D:118:ARG:HG2	2.07	0.54
1:A:200:LEU:O	1:A:201:GLU:CB	2.39	0.54
1:A:296:GLU:HG3	3:A:2218:HOH:O	2.07	0.54
1:B:103:ARG:CB	1:B:146:MSE:HE1	2.37	0.54
1:D:310:ASN:C	1:D:312:PRO:CD	2.76	0.54
1:A:201:GLU:O	1:A:202:LEU:HD12	2.08	0.53
1:D:13:LEU:HA	3:D:2006:HOH:O	2.08	0.53
1:D:313:MSE:CG	3:D:2048:HOH:O	2.14	0.53
1:B:97:ARG:HD2	2:B:401:ACO:H10	1.90	0.53
1:B:169:LEU:HD13	1:B:286:LEU:HD22	1.89	0.53
1:D:202:LEU:C	3:D:2207:HOH:O	2.52	0.53
1:A:202:LEU:HD13	3:A:2179:HOH:O	2.08	0.53
1:B:37:ARG:HD2	3:B:2042:HOH:O	2.08	0.53
1:B:84:THR:O	1:B:118:ARG:HG2	2.09	0.53
1:C:88:ASP:HB3	3:C:2091:HOH:O	2.07	0.52
1:C:97:ARG:HB3	2:C:401:ACO:O4A	2.09	0.52
1:A:195:VAL:HB	1:A:196:PRO:HD3	1.90	0.52
1:B:300:ARG:HH21	1:B:300:ARG:CG	2.23	0.52
1:D:276:ARG:CZ	3:D:2266:HOH:O	2.41	0.52
1:D:43:ASN:CG	3:D:2053:HOH:O	2.52	0.52
1:D:86:ARG:HD2	1:D:88:ASP:OD2	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:311:HIS:N	1:C:311:HIS:CD2	2.77	0.52
1:C:219:SER:O	1:C:223:GLN:HG3	2.10	0.52
1:B:296:GLU:HG3	3:B:2245:HOH:O	2.09	0.51
1:D:36:PRO:C	3:D:2034:HOH:O	2.38	0.51
1:A:106:TRP:CD1	1:A:146:MSE:CE	2.92	0.51
1:B:300:ARG:HG3	1:B:300:ARG:HH21	1.75	0.51
1:C:327:ARG:HD3	3:C:2265:HOH:O	2.09	0.51
1:D:129:LEU:CD1	1:D:327:ARG:HB2	2.40	0.51
1:B:310:ASN:ND2	1:B:310:ASN:C	2.69	0.51
1:B:327:ARG:HD3	3:B:2126:HOH:O	2.10	0.51
1:D:214:GLN:HG2	3:D:2204:HOH:O	2.11	0.51
1:A:151:SER:OG	1:A:152:ASP:N	2.44	0.51
1:B:110:ARG:NH1	3:B:2108:HOH:O	2.42	0.51
1:D:311:HIS:CD2	3:D:2288:HOH:O	2.65	0.50
1:B:211:ARG:N	1:B:211:ARG:NE	2.60	0.50
1:B:17:ASP:O	1:B:21:VAL:HG23	2.12	0.50
1:C:150:ARG:NH1	1:C:331:TRP:HZ2	2.10	0.50
1:A:63:VAL:CG1	1:A:70:VAL:HG22	2.42	0.50
1:B:202:LEU:HB2	1:B:270:VAL:HG13	1.93	0.50
1:D:43:ASN:HB2	3:D:2053:HOH:O	2.12	0.50
1:D:43:ASN:CB	3:D:2053:HOH:O	2.60	0.50
1:A:9:GLU:HB3	1:A:65:ARG:HG3	1.94	0.49
1:D:61:ASP:CG	1:D:75:ARG:HE	2.20	0.49
1:B:106:TRP:NE1	1:B:110:ARG:HD2	2.27	0.49
1:B:237:HIS:CD2	1:B:253:SER:HB3	2.47	0.49
1:B:312:PRO:O	1:B:316:VAL:HG23	2.13	0.49
1:D:173:VAL:HG23	3:D:2276:HOH:O	2.12	0.49
1:D:24:TRP:CH2	1:D:46:MSE:HE3	2.48	0.49
1:D:192:GLU:HG2	3:D:2198:HOH:O	2.13	0.49
1:B:202:LEU:CB	1:B:270:VAL:HG13	2.43	0.49
1:C:310:ASN:ND2	1:C:313:MSE:SE	2.95	0.48
1:A:103:ARG:HG3	1:A:146:MSE:HE1	1.94	0.48
1:B:334:GLU:HG3	3:B:2265:HOH:O	2.12	0.48
1:C:127:GLU:CD	1:C:150:ARG:NH1	2.71	0.48
1:C:150:ARG:NH1	1:C:331:TRP:CZ2	2.81	0.48
1:C:310:ASN:OD1	1:C:310:ASN:O	2.32	0.48
1:B:84:THR:C	1:B:118:ARG:HG2	2.38	0.48
1:B:129:LEU:O	1:B:130:PRO:C	2.57	0.48
1:D:106:TRP:NE1	1:D:110:ARG:HD2	2.28	0.48
1:C:234:ARG:HH21	1:C:296:GLU:CD	2.20	0.48
1:D:218:THR:CG2	3:D:2216:HOH:O	2.53	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:ARG:O	1:A:98:ARG:HB2	2.13	0.48
1:C:310:ASN:OD1	1:C:313:MSE:SE	2.81	0.48
1:D:201:GLU:O	1:D:202:LEU:CD1	2.44	0.48
1:B:320:LEU:CD1	1:B:320:LEU:N	2.77	0.47
1:D:256:LYS:HD2	1:D:296:GLU:OE1	2.14	0.47
1:A:121:LEU:HD12	1:A:121:LEU:N	2.29	0.47
1:A:151:SER:OG	1:A:153:ILE:HG12	2.14	0.47
1:D:84:THR:C	1:D:118:ARG:HG2	2.40	0.47
1:D:181:SER:HA	3:D:2184:HOH:O	2.14	0.47
1:A:135:GLN:O	1:A:136:ASP:C	2.57	0.47
1:A:312:PRO:O	1:A:316:VAL:HG23	2.13	0.47
1:A:129:LEU:HD13	1:A:327:ARG:HB2	1.95	0.47
1:D:297:PRO:HD2	3:D:2279:HOH:O	2.15	0.47
1:C:116:HIS:O	1:C:117:ASP:HB2	2.15	0.47
1:A:202:LEU:CD1	3:A:2179:HOH:O	2.63	0.46
1:A:234:ARG:HD3	3:A:2201:HOH:O	2.07	0.46
1:B:331:TRP:C	3:B:2261:HOH:O	2.53	0.46
1:C:103:ARG:HA	1:C:146:MSE:HE2	1.97	0.46
1:D:311:HIS:N	1:D:312:PRO:HD3	2.30	0.46
1:A:266:GLN:HG2	3:A:2226:HOH:O	2.14	0.46
1:B:62:TRP:CD2	1:B:108:HIS:HE1	2.33	0.46
1:C:99:ARG:NE	3:C:2108:HOH:O	2.48	0.46
1:D:11:VAL:HG13	1:D:12:PRO:HD2	1.97	0.46
1:C:114:ARG:HG2	3:C:2274:HOH:O	2.15	0.46
1:A:24:TRP:HE1	1:A:89:GLN:NE2	2.04	0.46
1:B:273:ARG:HG3	3:B:2228:HOH:O	2.15	0.46
1:D:145:ALA:C	1:D:146:MSE:HE2	2.41	0.45
1:A:153:ILE:HG12	1:A:153:ILE:H	1.47	0.45
1:C:311:HIS:CD2	1:C:311:HIS:H	2.34	0.45
1:C:218:THR:HG22	3:C:2185:HOH:O	2.17	0.45
1:D:316:VAL:CG1	3:D:2273:HOH:O	2.44	0.45
1:B:200:LEU:HA	1:B:200:LEU:HD12	1.67	0.45
1:D:84:THR:HG22	1:D:85:ALA:N	2.31	0.45
1:D:234:ARG:CD	3:D:2234:HOH:O	2.36	0.45
1:A:150:ARG:NH1	1:A:331:TRP:CZ2	2.85	0.45
1:A:200:LEU:HD11	1:A:272:HIS:HB2	1.98	0.45
1:B:84:THR:HG22	1:B:85:ALA:N	2.31	0.45
1:C:116:HIS:O	1:C:118:ARG:HD2	2.17	0.45
1:C:316:VAL:HG12	3:C:2236:HOH:O	2.02	0.45
1:B:131:SER:OG	1:B:132:GLY:N	2.49	0.45
1:A:98:ARG:NH2	3:A:2098:HOH:O	2.07	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:ARG:NH1	1:A:331:TRP:HZ2	2.15	0.44
1:A:195:VAL:N	1:A:196:PRO:CD	2.80	0.44
1:D:164:ASP:OD1	1:D:300:ARG:HD2	2.18	0.44
1:C:64:VAL:HG13	1:C:71:VAL:HG23	2.00	0.44
1:D:62:TRP:CD2	1:D:108:HIS:HE1	2.34	0.44
1:C:24:TRP:CZ2	1:C:46:MSE:HE2	2.52	0.44
1:C:256:LYS:HD3	1:C:257:THR:N	2.33	0.44
1:D:42:CYS:O	1:D:45:ASP:HB3	2.18	0.44
1:D:187:THR:HG21	3:D:2191:HOH:O	2.18	0.44
1:A:285:LYS:CE	3:A:2239:HOH:O	2.65	0.44
1:B:300:ARG:NH1	3:B:2249:HOH:O	2.50	0.44
1:D:200:LEU:O	1:D:201:GLU:CB	2.55	0.44
1:D:47:VAL:O	1:D:48:GLY:C	2.59	0.44
1:B:152:ASP:C	1:B:152:ASP:OD1	2.61	0.43
1:A:21:VAL:O	1:A:25:LEU:HG	2.18	0.43
1:B:28:MSE:HE1	1:B:46:MSE:HG3	2.00	0.43
1:C:129:LEU:O	1:C:130:PRO:C	2.62	0.43
1:A:311:HIS:HB2	1:A:312:PRO:HD3	1.99	0.43
1:C:121:LEU:HD12	1:C:121:LEU:N	2.33	0.43
1:C:237:HIS:CD2	1:C:253:SER:HB3	2.53	0.43
1:D:97:ARG:HB3	2:D:401:ACO:O4A	2.18	0.43
1:B:158:HIS:HB3	1:B:326:ASP:OD1	2.18	0.43
1:C:192:GLU:N	3:C:2174:HOH:O	2.31	0.43
1:D:62:TRP:CE3	1:D:108:HIS:HE1	2.36	0.43
1:D:313:MSE:O	1:D:314:ILE:C	2.61	0.43
1:A:120:THR:HA	1:A:333:ALA:O	2.18	0.43
1:B:37:ARG:NH1	3:B:2042:HOH:O	2.36	0.43
1:B:280:LEU:O	1:B:284:LEU:CD2	2.67	0.43
1:A:237:HIS:CD2	1:A:253:SER:HB3	2.54	0.43
1:B:150:ARG:HD3	1:B:329:VAL:HG11	2.00	0.43
1:A:211:ARG:HD2	1:A:213:ALA:HB2	1.99	0.43
1:B:20:THR:O	1:B:21:VAL:C	2.62	0.43
1:B:202:LEU:HD21	1:B:276:ARG:HH21	1.82	0.42
1:B:129:LEU:N	3:B:2127:HOH:O	2.38	0.42
1:B:92:VAL:O	1:B:97:ARG:NH1	2.43	0.42
1:D:92:VAL:O	1:D:93:HIS:C	2.62	0.42
1:D:169:LEU:HD22	1:D:289:LEU:HD23	2.01	0.42
1:B:199:GLU:O	1:B:201:GLU:N	2.53	0.42
1:D:313:MSE:CB	3:D:2048:HOH:O	2.57	0.42
1:A:152:ASP:OD1	1:A:152:ASP:C	2.62	0.42
1:A:169:LEU:HD13	1:A:286:LEU:HD22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:286:LEU:HD23	1:D:286:LEU:HA	1.77	0.42
1:B:272:HIS:ND1	1:B:274:GLU:N	2.51	0.42
1:B:317:ASN:HD22	1:B:317:ASN:HA	1.70	0.42
1:C:234:ARG:CD	3:C:2200:HOH:O	2.62	0.41
1:A:16:ASN:OD1	1:A:16:ASN:C	2.62	0.41
1:C:298:GLU:HB2	3:C:2243:HOH:O	2.20	0.41
1:B:121:LEU:N	1:B:121:LEU:HD12	2.35	0.41
1:D:43:ASN:ND2	3:D:2053:HOH:O	2.52	0.41
1:A:251:TYR:N	1:A:251:TYR:CD1	2.88	0.41
2:D:401:ACO:P2A	2:D:401:ACO:H133	2.56	0.41
1:B:93:HIS:CD2	1:B:95:GLY:N	2.63	0.41
1:B:200:LEU:HD23	1:B:251:TYR:CE1	2.56	0.41
1:C:106:TRP:HD1	1:C:146:MSE:HE3	1.79	0.41
1:A:93:HIS:CD2	1:A:95:GLY:H	2.21	0.41
1:A:90:LEU:HD23	1:A:90:LEU:C	2.45	0.41
1:A:316:VAL:HG11	3:A:2239:HOH:O	2.20	0.41
1:C:158:HIS:HB3	1:C:326:ASP:OD1	2.21	0.41
1:C:65:ARG:HA	1:C:69:ARG:O	2.20	0.41
1:C:86:ARG:HD2	1:C:88:ASP:OD2	2.21	0.41
1:D:84:THR:HG22	1:D:85:ALA:H	1.86	0.41
1:A:86:ARG:HG3	3:A:2119:HOH:O	2.20	0.41
1:A:286:LEU:HA	1:A:286:LEU:HD23	1.74	0.41
1:B:256:LYS:HD2	1:B:296:GLU:CD	2.45	0.41
1:D:158:HIS:HB3	1:D:326:ASP:OD1	2.21	0.41
1:D:65:ARG:NH1	3:D:2076:HOH:O	2.53	0.40
1:B:106:TRP:CB	1:B:146:MSE:HE3	2.41	0.40
1:B:223:GLN:NE2	3:B:2194:HOH:O	2.54	0.40
1:D:256:LYS:HE2	1:D:263:TYR:O	2.21	0.40
1:A:110:ARG:HH11	1:A:110:ARG:HD3	1.76	0.40
1:A:150:ARG:HD3	1:A:329:VAL:HG11	2.04	0.40
1:C:61:ASP:CG	1:C:75:ARG:HE	2.30	0.40
1:C:254:VAL:HG23	3:C:2237:HOH:O	2.22	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)
2:C:401:ACO:O7A	2:D:401:ACO:O2B[2_555]	2.11	0.09

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	324/339 (96%)	315 (97%)	8 (2%)	1 (0%)	36	48
1	B	322/339 (95%)	312 (97%)	8 (2%)	2 (1%)	21	30
1	C	321/339 (95%)	312 (97%)	8 (2%)	1 (0%)	36	48
1	D	321/339 (95%)	305 (95%)	13 (4%)	3 (1%)	14	20
All	All	1288/1356 (95%)	1244 (97%)	37 (3%)	7 (0%)	24	34

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	201	GLU
1	B	213	ALA
1	D	201	GLU
1	B	130	PRO
1	D	131	SER
1	D	132	GLY
1	C	312	PRO

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	248/249 (100%)	224 (90%)	24 (10%)	8	11
1	B	247/249 (99%)	228 (92%)	19 (8%)	12	18
1	C	247/249 (99%)	226 (92%)	21 (8%)	10	15

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	247/249 (99%)	231 (94%)	16 (6%)	15	24
All	All	989/996 (99%)	909 (92%)	80 (8%)	11	16

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	THR
1	A	37	ARG
1	A	63	VAL
1	A	64	VAL
1	A	65	ARG
1	A	78	LEU
1	A	97	ARG
1	A	114	ARG
1	A	129	LEU
1	A	146	MSE
1	A	153	ILE
1	A	163	LEU
1	A	165	ARG
1	A	169	LEU
1	A	215	GLU
1	A	216	VAL
1	A	256	LYS
1	A	268	MSE
1	A	280	LEU
1	A	284	LEU
1	A	293	LEU
1	A	296	GLU
1	A	300	ARG
1	A	320	LEU
1	B	37	ARG
1	B	63	VAL
1	B	65	ARG
1	B	97	ARG
1	B	130	PRO
1	B	163	LEU
1	B	169	LEU
1	B	200	LEU
1	B	201	GLU
1	B	211	ARG
1	B	256	LYS
1	B	268	MSE

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Mol	Chain	Res	Type
1	B	280	LEU
1	B	284	LEU
1	B	293	LEU
1	B	300	ARG
1	B	310	ASN
1	B	313	MSE
1	B	320	LEU
1	C	7	GLU
1	C	37	ARG
1	C	63	VAL
1	C	97	ARG
1	C	118	ARG
1	C	129	LEU
1	C	131	SER
1	C	146	MSE
1	C	163	LEU
1	C	169	LEU
1	C	200	LEU
1	C	202	LEU
1	C	211	ARG
1	C	218	THR
1	C	256	LYS
1	C	268	MSE
1	C	280	LEU
1	C	284	LEU
1	C	293	LEU
1	C	300	ARG
1	C	320	LEU
1	D	37	ARG
1	D	63	VAL
1	D	78	LEU
1	D	97	ARG
1	D	163	LEU
1	D	169	LEU
1	D	202	LEU
1	D	215	GLU
1	D	218	THR
1	D	256	LYS
1	D	276	ARG
1	D	280	LEU
1	D	284	LEU
1	D	293	LEU

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Mol	Chain	Res	Type
1	D	296	GLU
1	D	320	LEU

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

Mol	Chain	Res	Type
1	A	89	GLN
1	A	93	HIS
1	A	108	HIS
1	A	214	GLN
1	A	278	HIS
1	A	306	ASN
1	A	317	ASN
1	B	89	GLN
1	B	93	HIS
1	B	108	HIS
1	B	159	GLN
1	B	278	HIS
1	B	306	ASN
1	B	317	ASN
1	C	89	GLN
1	C	108	HIS
1	C	214	GLN
1	C	288	ASN
1	C	310	ASN
1	C	311	HIS
1	D	89	GLN
1	D	93	HIS
1	D	108	HIS
1	D	159	GLN
1	D	214	GLN
1	D	223	GLN
1	D	260	ASN
1	D	278	HIS
1	D	306	ASN
1	D	311	HIS
1	D	317	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	ACO	B	401	-	51,53,53	1.80	11 (21%)	73,79,79	3.37	27 (36%)
2	ACO	A	401	-	51,53,53	1.71	11 (21%)	73,79,79	3.56	27 (36%)
2	ACO	D	401	-	51,53,53	1.89	9 (17%)	73,79,79	3.51	28 (38%)
2	ACO	C	401	-	51,53,53	1.56	7 (13%)	73,79,79	3.58	28 (38%)

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	ACO	B	401	-	-	11/51/67/67	0/3/3/3
2	ACO	A	401	-	-	15/51/67/67	0/3/3/3
2	ACO	D	401	-	-	15/51/67/67	0/3/3/3
2	ACO	C	401	-	-	14/51/67/67	0/3/3/3

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	401	ACO	P3B-O7A	6.01	1.69	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	ACO	CDP-CBP	-5.19	1.43	1.53
2	B	401	ACO	P3B-O7A	4.96	1.65	1.50
2	C	401	ACO	P3B-O7A	4.65	1.64	1.50
2	D	401	ACO	P3B-O3B	4.50	1.67	1.59
2	D	401	ACO	P2A-O4A	4.45	1.66	1.50
2	D	401	ACO	CDP-CBP	-4.34	1.44	1.53
2	A	401	ACO	P3B-O7A	4.16	1.63	1.50
2	A	401	ACO	P2A-O3A	4.12	1.64	1.59
2	D	401	ACO	O4B-C1B	3.81	1.50	1.42
2	C	401	ACO	P2A-O4A	3.74	1.63	1.50
2	A	401	ACO	CCP-CBP	3.71	1.58	1.52
2	B	401	ACO	C5A-N7A	-3.69	1.32	1.39
2	A	401	ACO	P3B-O3B	3.54	1.65	1.59
2	A	401	ACO	P2A-O4A	3.45	1.62	1.50
2	B	401	ACO	O4B-C1B	3.37	1.49	1.42
2	B	401	ACO	P2A-O4A	3.19	1.61	1.50
2	D	401	ACO	P2A-O6A	3.19	1.71	1.59
2	B	401	ACO	C1B-N9A	3.17	1.55	1.46
2	C	401	ACO	C5A-N7A	-3.17	1.33	1.39
2	C	401	ACO	O6A-CCP	-3.08	1.34	1.43
2	A	401	ACO	O5B-C5B	-2.96	1.33	1.44
2	D	401	ACO	C5A-N7A	-2.93	1.33	1.39
2	C	401	ACO	O5B-C5B	-2.90	1.33	1.44
2	C	401	ACO	O4B-C1B	2.88	1.48	1.42
2	C	401	ACO	P1A-O1A	2.81	1.60	1.50
2	B	401	ACO	O5B-C5B	-2.79	1.34	1.44
2	A	401	ACO	P1A-O3A	-2.74	1.56	1.59
2	B	401	ACO	CCP-CBP	2.62	1.56	1.52
2	D	401	ACO	P2A-O3A	-2.49	1.56	1.59
2	D	401	ACO	P1A-O1A	2.49	1.59	1.50
2	A	401	ACO	O6A-CCP	-2.45	1.36	1.43
2	A	401	ACO	C5A-N7A	-2.43	1.34	1.39
2	B	401	ACO	C7P-N8P	2.28	1.51	1.46
2	B	401	ACO	P2A-O6A	2.25	1.68	1.59
2	A	401	ACO	C4A-N9A	-2.23	1.33	1.37
2	A	401	ACO	P2A-O6A	2.16	1.67	1.59
2	B	401	ACO	P3B-O3B	-2.14	1.55	1.59

All (110) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	ACO	CEP-CBP-CCP	-15.18	83.17	108.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	ACO	CDP-CBP-CCP	-14.40	84.46	108.22
2	C	401	ACO	CEP-CBP-CCP	-14.29	84.63	108.22
2	C	401	ACO	CDP-CBP-CCP	-12.83	87.04	108.22
2	B	401	ACO	CDP-CBP-CCP	-12.42	87.73	108.22
2	A	401	ACO	CDP-CBP-CCP	-12.30	87.91	108.22
2	D	401	ACO	CEP-CBP-CCP	-10.75	90.48	108.22
2	B	401	ACO	CEP-CBP-CCP	-9.55	92.46	108.22
2	D	401	ACO	O9A-P3B-O8A	-8.49	75.97	107.80
2	B	401	ACO	P3B-O3B-C3B	-8.28	101.33	123.43
2	C	401	ACO	O9A-P3B-O8A	-8.28	76.77	107.80
2	A	401	ACO	O9A-P3B-O8A	-8.24	76.91	107.80
2	B	401	ACO	CEP-CBP-CAP	7.56	121.66	108.77
2	B	401	ACO	CDP-CBP-CAP	7.51	121.57	108.77
2	A	401	ACO	CEP-CBP-CAP	7.34	121.28	108.77
2	A	401	ACO	O9A-P3B-O3B	-7.32	77.33	105.85
2	C	401	ACO	CEP-CBP-CAP	7.27	121.17	108.77
2	D	401	ACO	O9A-P3B-O3B	-7.26	77.57	105.85
2	D	401	ACO	CDP-CBP-CAP	7.24	121.12	108.77
2	C	401	ACO	CEP-CBP-CDP	6.69	122.53	109.20
2	C	401	ACO	O9A-P3B-O3B	-6.56	80.30	105.85
2	D	401	ACO	O9A-P3B-O7A	-6.53	85.39	110.83
2	C	401	ACO	O9A-P3B-O7A	-6.51	85.48	110.83
2	A	401	ACO	O9A-P3B-O7A	-6.32	86.21	110.83
2	A	401	ACO	CEP-CBP-CDP	6.32	121.80	109.20
2	D	401	ACO	N3A-C2A-N1A	-6.27	119.09	128.58
2	C	401	ACO	N3A-C2A-N1A	-6.25	119.11	128.58
2	D	401	ACO	C5A-C4A-N3A	-6.17	118.22	126.72
2	D	401	ACO	CEP-CBP-CDP	6.02	121.21	109.20
2	B	401	ACO	O8A-P3B-O7A	5.63	132.75	110.83
2	A	401	ACO	C5A-C4A-N3A	-5.61	118.99	126.72
2	B	401	ACO	C5A-C4A-N3A	-5.57	119.05	126.72
2	B	401	ACO	O9A-P3B-O7A	-5.56	89.16	110.83
2	B	401	ACO	N3A-C2A-N1A	-5.54	120.19	128.58
2	C	401	ACO	C5B-C4B-C3B	-5.38	96.46	114.38
2	C	401	ACO	C5A-C4A-N3A	-5.37	119.33	126.72
2	B	401	ACO	O9A-P3B-O8A	-5.33	87.80	107.80
2	D	401	ACO	CEP-CBP-CAP	5.19	117.61	108.77
2	B	401	ACO	O4B-C1B-N9A	5.08	117.84	108.09
2	A	401	ACO	N3A-C2A-N1A	-4.94	121.11	128.58
2	D	401	ACO	C2A-N3A-C4A	4.83	123.63	111.83
2	B	401	ACO	C2A-N3A-C4A	4.39	122.54	111.83
2	C	401	ACO	C2A-N3A-C4A	4.37	122.50	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	ACO	O9A-P3B-O3B	-4.34	88.94	105.85
2	A	401	ACO	CDP-CBP-CAP	4.32	116.14	108.77
2	B	401	ACO	N9A-C8A-N7A	-4.25	107.91	113.94
2	C	401	ACO	N3A-C4A-N9A	4.12	134.18	127.17
2	A	401	ACO	C5B-C4B-C3B	-4.12	100.66	114.38
2	B	401	ACO	C5A-N7A-C8A	4.11	109.92	103.45
2	D	401	ACO	N3A-C4A-N9A	4.10	134.14	127.17
2	A	401	ACO	C2A-N3A-C4A	4.10	121.83	111.83
2	B	401	ACO	O8A-P3B-O3B	4.05	121.65	105.85
2	C	401	ACO	CDP-CBP-CAP	3.93	115.46	108.77
2	B	401	ACO	CEP-CBP-CDP	3.74	116.66	109.20
2	B	401	ACO	C4A-C5A-N7A	-3.69	106.36	110.58
2	D	401	ACO	O4B-C4B-C5B	-3.69	97.51	109.33
2	A	401	ACO	N9A-C8A-N7A	-3.69	108.70	113.94
2	C	401	ACO	O2A-P1A-O5B	3.68	124.26	107.57
2	D	401	ACO	N9A-C8A-N7A	-3.61	108.82	113.94
2	A	401	ACO	N3A-C4A-N9A	3.55	133.21	127.17
2	B	401	ACO	O4B-C4B-C5B	-3.40	98.44	109.33
2	C	401	ACO	C6P-C7P-N8P	-3.33	104.91	112.00
2	C	401	ACO	O3A-P2A-O4A	-3.30	100.79	110.70
2	A	401	ACO	C4A-C5A-N7A	-3.28	106.83	110.58
2	A	401	ACO	O4B-C1B-N9A	3.28	114.39	108.09
2	D	401	ACO	O5B-P1A-O1A	-3.23	96.14	108.94
2	C	401	ACO	O8A-P3B-O3B	3.18	118.23	105.85
2	A	401	ACO	C5A-N7A-C8A	3.12	108.35	103.45
2	D	401	ACO	O8A-P3B-O3B	3.09	117.90	105.85
2	B	401	ACO	N3A-C4A-N9A	3.09	132.42	127.17
2	D	401	ACO	C5A-N7A-C8A	3.02	108.20	103.45
2	B	401	ACO	O5A-P2A-O6A	2.99	121.13	107.57
2	A	401	ACO	O4B-C4B-C5B	-2.99	99.77	109.33
2	B	401	ACO	C6P-C7P-N8P	-2.95	105.73	112.00
2	A	401	ACO	O5A-P2A-O3A	2.92	115.16	107.27
2	C	401	ACO	C2B-C3B-C4B	2.88	108.28	103.24
2	C	401	ACO	C2B-C1B-N9A	-2.85	106.22	113.30
2	A	401	ACO	O2A-P1A-O5B	2.84	120.43	107.57
2	D	401	ACO	C2B-C3B-C4B	2.81	108.16	103.24
2	B	401	ACO	C2B-C3B-C4B	2.79	108.12	103.24
2	A	401	ACO	P3B-O3B-C3B	-2.76	116.05	123.43
2	D	401	ACO	O3A-P2A-O4A	-2.76	102.39	110.70
2	A	401	ACO	O3A-P2A-O4A	-2.71	102.55	110.70
2	C	401	ACO	N9A-C8A-N7A	-2.71	110.10	113.94
2	D	401	ACO	O2A-P1A-O5B	2.70	119.81	107.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	ACO	O6A-CCP-CBP	2.69	114.87	110.55
2	A	401	ACO	O8A-P3B-O3B	2.68	116.28	105.85
2	D	401	ACO	C4A-C5A-N7A	-2.67	107.53	110.58
2	D	401	ACO	P3B-O3B-C3B	-2.62	116.44	123.43
2	D	401	ACO	O4B-C1B-N9A	2.62	113.11	108.09
2	D	401	ACO	OAP-CAP-CBP	-2.60	104.15	110.18
2	C	401	ACO	O5A-P2A-O3A	2.56	114.18	107.27
2	B	401	ACO	O6A-CCP-CBP	2.52	114.59	110.55
2	B	401	ACO	C5A-C4A-N9A	2.48	108.51	105.81
2	C	401	ACO	O5B-P1A-O1A	-2.46	99.17	108.94
2	A	401	ACO	C4A-N9A-C8A	2.46	108.32	105.74
2	C	401	ACO	P3B-O3B-C3B	-2.43	116.94	123.43
2	C	401	ACO	CAP-C9P-N8P	-2.43	111.88	116.48
2	D	401	ACO	C5B-C4B-C3B	-2.36	106.51	114.38
2	A	401	ACO	O3B-P3B-O7A	2.35	117.71	109.33
2	C	401	ACO	C5A-N7A-C8A	2.29	107.04	103.45
2	A	401	ACO	O6A-CCP-CBP	2.28	114.22	110.55
2	A	401	ACO	OAP-CAP-CBP	-2.28	104.91	110.18
2	B	401	ACO	C6A-C5A-N7A	2.27	136.46	132.09
2	B	401	ACO	OAP-CAP-CBP	-2.21	105.07	110.18
2	D	401	ACO	C2P-C3P-N4P	-2.11	108.00	112.41
2	C	401	ACO	O3B-P3B-O7A	2.11	116.85	109.33
2	D	401	ACO	C7P-C6P-C5P	-2.10	108.89	112.39
2	C	401	ACO	C3P-N4P-C5P	-2.08	118.95	122.82
2	C	401	ACO	O4B-C1B-N9A	2.04	112.01	108.09

There are no chirality outliers.

All (55) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	ACO	CDP-CBP-CCP-O6A
2	A	401	ACO	CEP-CBP-CCP-O6A
2	A	401	ACO	CAP-CBP-CCP-O6A
2	A	401	ACO	OAP-CAP-CBP-CCP
2	A	401	ACO	C9P-CAP-CBP-CCP
2	A	401	ACO	C9P-CAP-CBP-CEP
2	B	401	ACO	C5B-O5B-P1A-O1A
2	B	401	ACO	CDP-CBP-CCP-O6A
2	B	401	ACO	CEP-CBP-CCP-O6A
2	B	401	ACO	CAP-CBP-CCP-O6A
2	B	401	ACO	OAP-CAP-CBP-CCP
2	B	401	ACO	C9P-CAP-CBP-CCP

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Mol	Chain	Res	Type	Atoms
2	B	401	ACO	OAP-CAP-CBP-CEP
2	B	401	ACO	C9P-CAP-CBP-CEP
2	C	401	ACO	CCP-O6A-P2A-O4A
2	C	401	ACO	CDP-CBP-CCP-O6A
2	C	401	ACO	CEP-CBP-CCP-O6A
2	C	401	ACO	CAP-CBP-CCP-O6A
2	C	401	ACO	OAP-CAP-CBP-CCP
2	C	401	ACO	C9P-CAP-CBP-CCP
2	C	401	ACO	C9P-CAP-CBP-CEP
2	D	401	ACO	C3B-C4B-C5B-O5B
2	D	401	ACO	O4B-C4B-C5B-O5B
2	D	401	ACO	C5B-O5B-P1A-O2A
2	D	401	ACO	C5B-O5B-P1A-O3A
2	D	401	ACO	CDP-CBP-CCP-O6A
2	D	401	ACO	CEP-CBP-CCP-O6A
2	D	401	ACO	CAP-CBP-CCP-O6A
2	D	401	ACO	OAP-CAP-CBP-CCP
2	D	401	ACO	C9P-CAP-CBP-CCP
2	D	401	ACO	OAP-CAP-CBP-CEP
2	D	401	ACO	C9P-CAP-CBP-CEP
2	A	401	ACO	C3B-C4B-C5B-O5B
2	A	401	ACO	O4B-C4B-C5B-O5B
2	A	401	ACO	OAP-CAP-CBP-CDP
2	A	401	ACO	OAP-CAP-CBP-CEP
2	C	401	ACO	OAP-CAP-CBP-CDP
2	C	401	ACO	OAP-CAP-CBP-CEP
2	A	401	ACO	C9P-CAP-CBP-CDP
2	C	401	ACO	O4B-C4B-C5B-O5B
2	C	401	ACO	P1A-O3A-P2A-O5A
2	D	401	ACO	P1A-O3A-P2A-O5A
2	C	401	ACO	C6P-C7P-N8P-C9P
2	A	401	ACO	CCP-O6A-P2A-O4A
2	B	401	ACO	C5B-O5B-P1A-O3A
2	A	401	ACO	P1A-O3A-P2A-O5A
2	C	401	ACO	C3B-O3B-P3B-O9A
2	D	401	ACO	C6P-C7P-N8P-C9P
2	A	401	ACO	P1A-O3A-P2A-O4A
2	B	401	ACO	P1A-O3A-P2A-O4A
2	C	401	ACO	P1A-O3A-P2A-O4A
2	A	401	ACO	C3B-O3B-P3B-O8A
2	D	401	ACO	CBP-CCP-O6A-P2A
2	D	401	ACO	C3P-C2P-S1P-C

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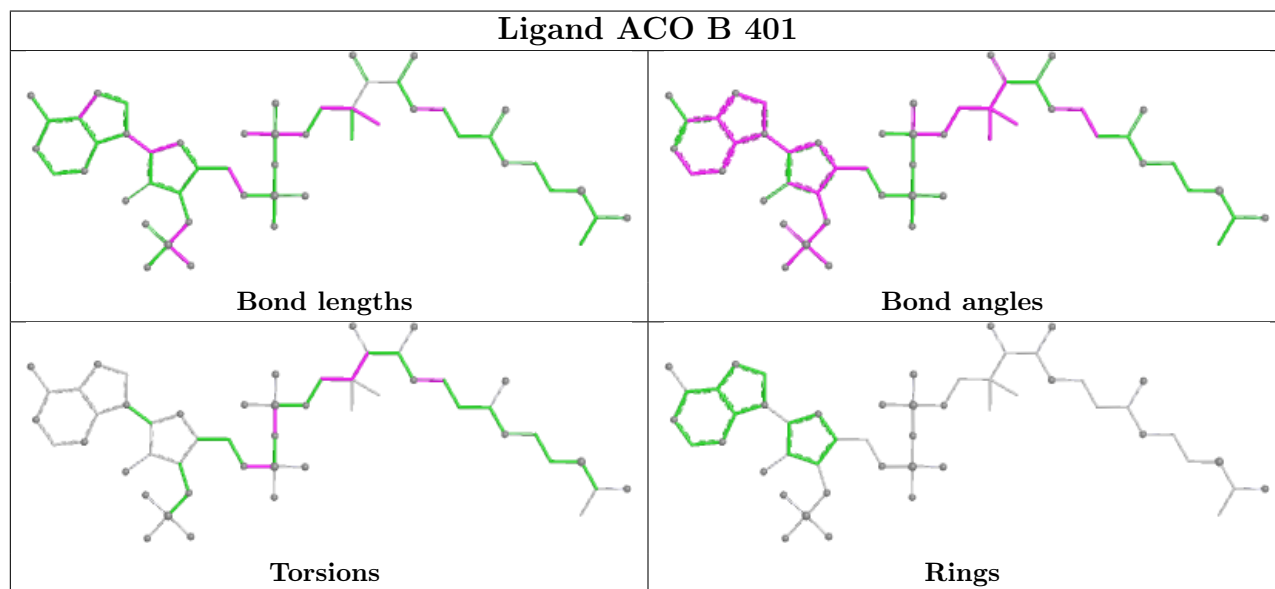
Mol	Chain	Res	Type	Atoms
2	B	401	ACO	C6P-C7P-N8P-C9P

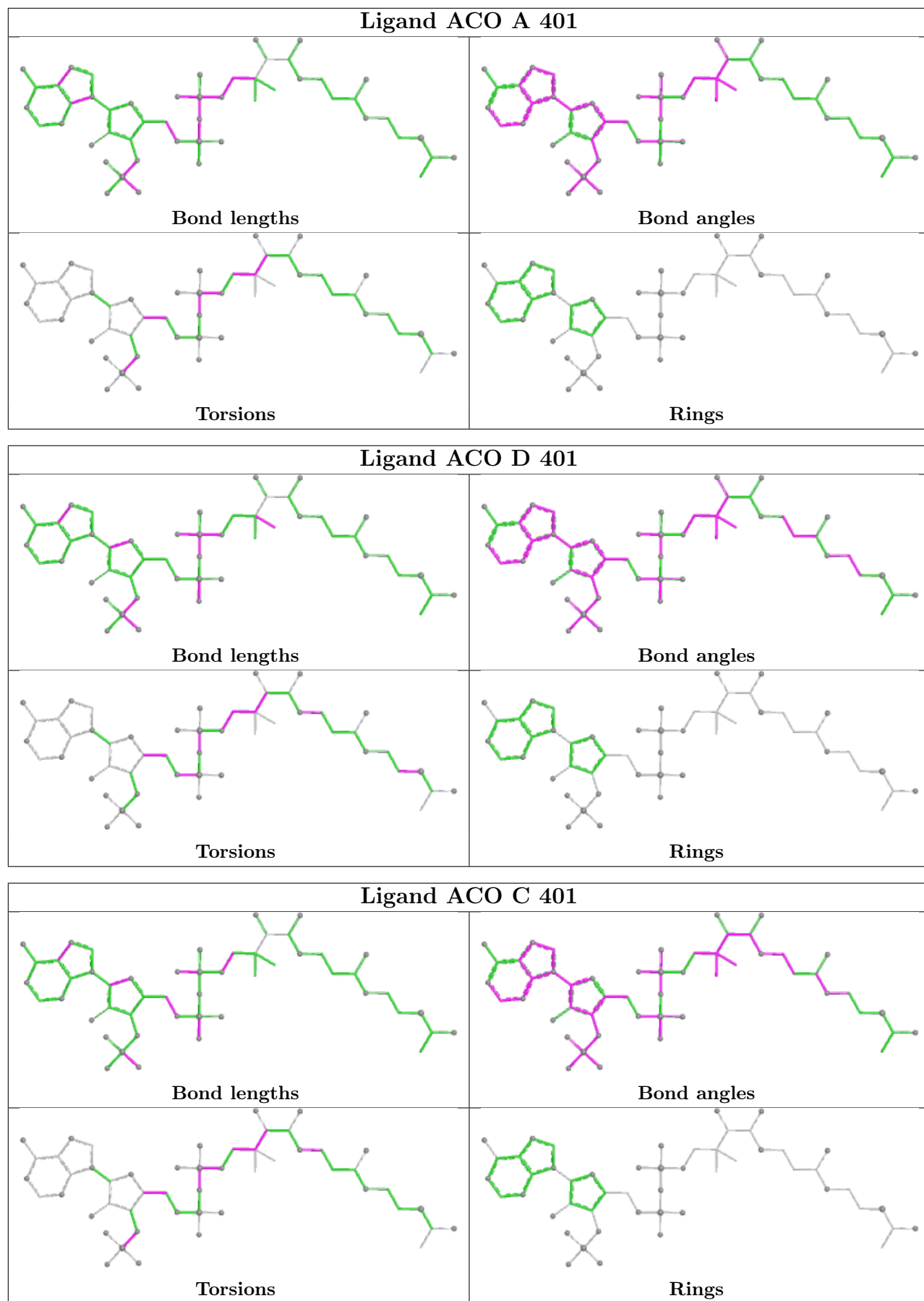
There are no ring outliers.

4 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	ACO	5	0
2	A	401	ACO	5	0
2	D	401	ACO	8	1
2	C	401	ACO	4	1

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

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	322/339 (94%)	-0.20	7 (2%) 62 61	15, 26, 53, 86	0
1	B	320/339 (94%)	-0.17	8 (2%) 58 58	13, 27, 55, 76	0
1	C	319/339 (94%)	-0.21	5 (1%) 70 69	16, 26, 51, 76	0
1	D	319/339 (94%)	-0.18	5 (1%) 70 69	14, 27, 52, 73	0
All	All	1280/1356 (94%)	-0.19	25 (1%) 65 63	13, 27, 52, 86	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	212	ALA	3.8
1	B	202	LEU	3.3
1	A	4	THR	3.1
1	B	213	ALA	3.1
1	D	212	ALA	3.0
1	A	5	ALA	3.0
1	C	212	ALA	3.0
1	A	213	ALA	2.9
1	B	6	GLY	2.9
1	C	213	ALA	2.8
1	A	202	LEU	2.8
1	D	213	ALA	2.6
1	B	212	ALA	2.6
1	B	312	PRO	2.6
1	C	312	PRO	2.5
1	C	202	LEU	2.4
1	D	312	PRO	2.3
1	C	200	LEU	2.3
1	B	211	ARG	2.2
1	D	310	ASN	2.2
1	A	6	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	314	ILE	2.1
1	D	131	SER	2.1
1	B	130	PRO	2.0
1	A	211	ARG	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

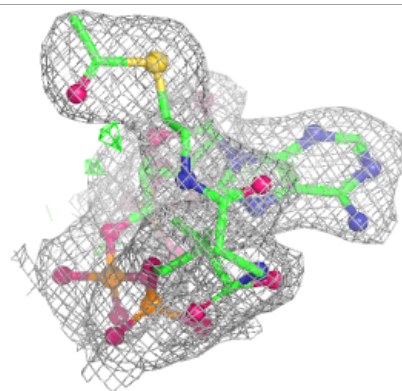
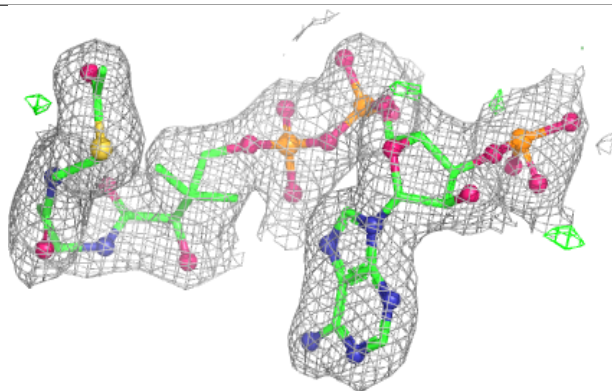
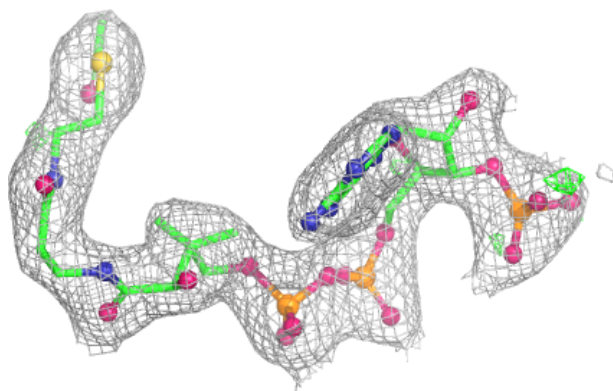
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q < 0.9
2	ACO	D	401	51/51	0.95	0.08	26,33,47,49	0
2	ACO	B	401	51/51	0.96	0.07	17,28,40,44	0
2	ACO	C	401	51/51	0.96	0.07	15,29,39,43	0
2	ACO	A	401	51/51	0.96	0.07	13,26,37,43	0

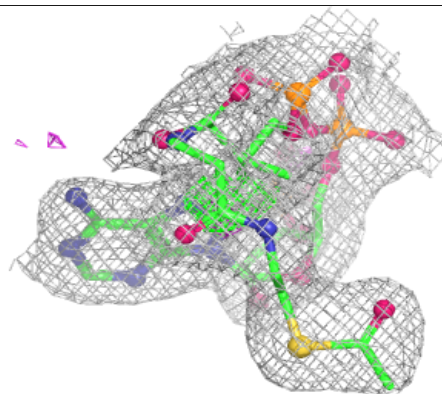
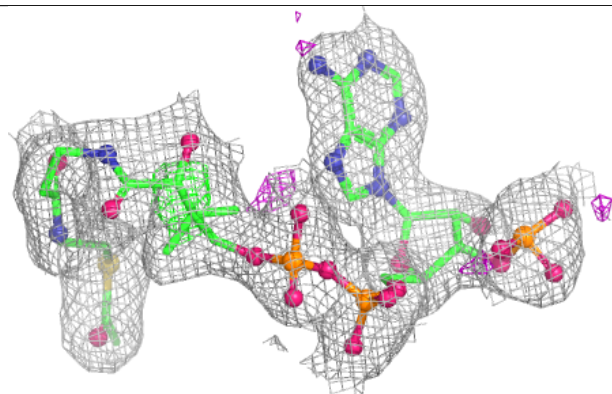
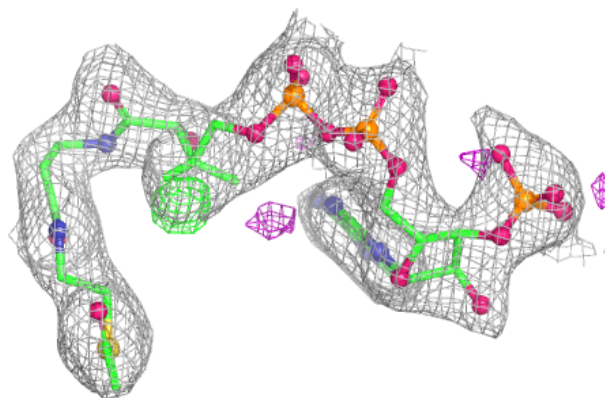
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ACO D 401:

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

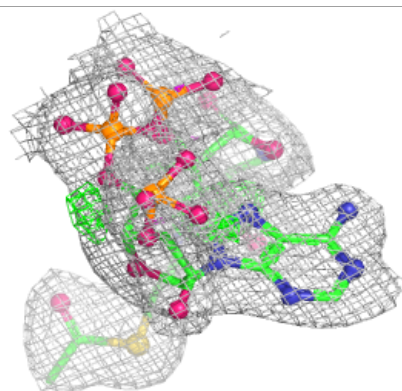
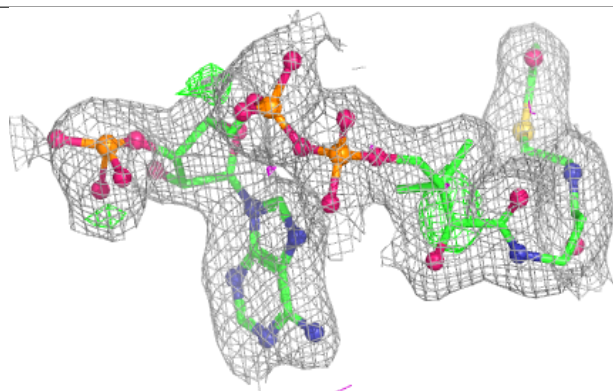
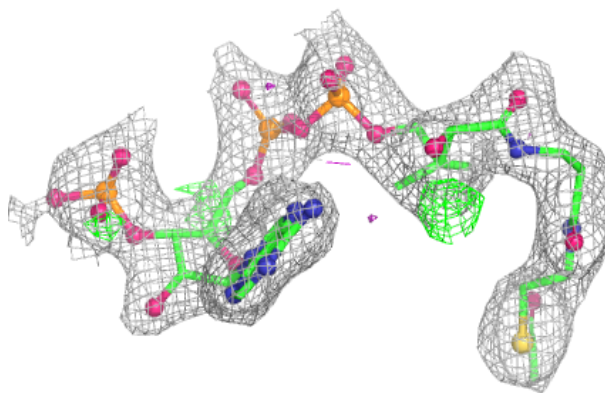
**Electron density around ACO B 401:**

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

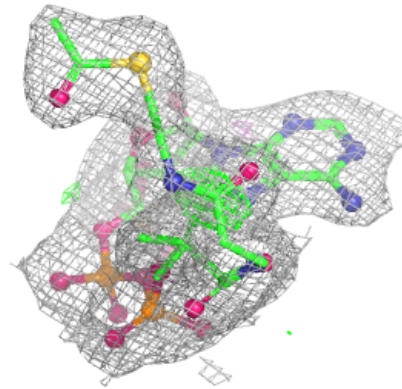
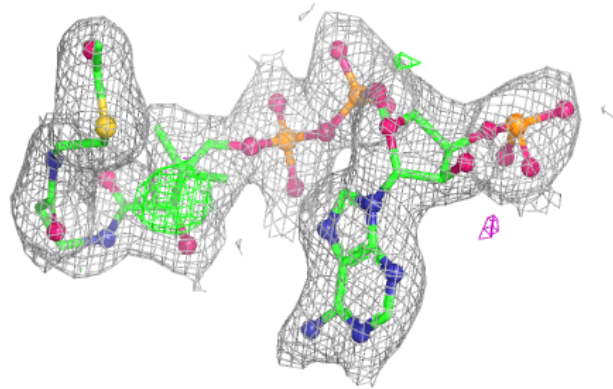
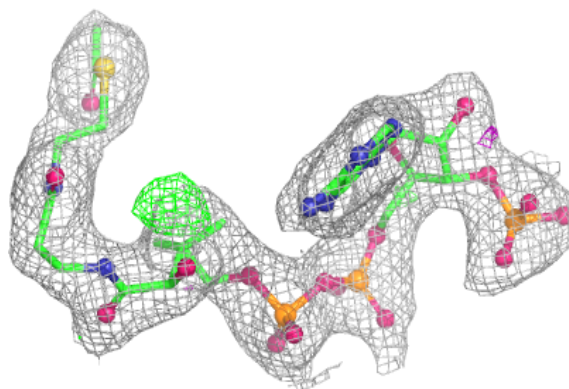


Electron density around ACO C 401:

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

**Electron density around ACO A 401:**

$2mF_o-DF_c$ (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.