



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 11:50 pm BST

PDB ID : 4AVC
Title : Crystal structure of protein lysine acetyltransferase Rv0998 in complex with acetyl CoA and cAMP
Authors : Lee, H.J.; Lang, P.T.; Fortune, S.M.; Sasseti, C.M.; Alber, T.
Deposited on : 2012-05-24
Resolution : 2.81 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 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 : 2.11
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) : 2.11

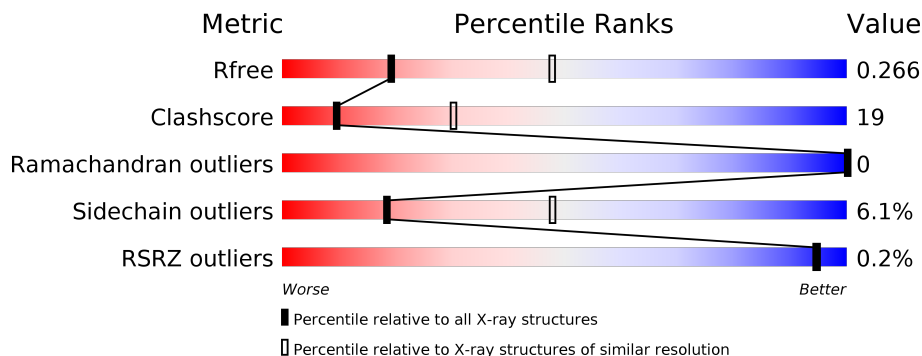
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on 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	333	
1	B	333	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ACO	A	1334	X	-	-	-
3	ACO	B	1334	X	-	-	-
4	PEG	A	1335	-	-	X	-

2 Entry composition [i](#)

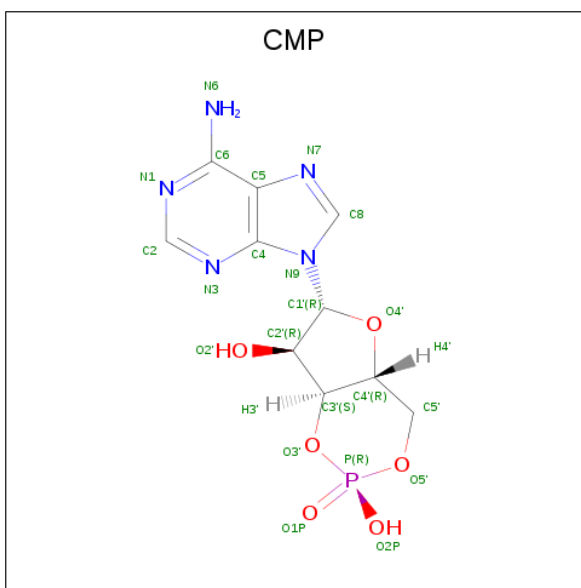
There are 6 unique types of molecules in this entry. The entry contains 5179 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called LYSINE ACETYLTRANSFERASE.

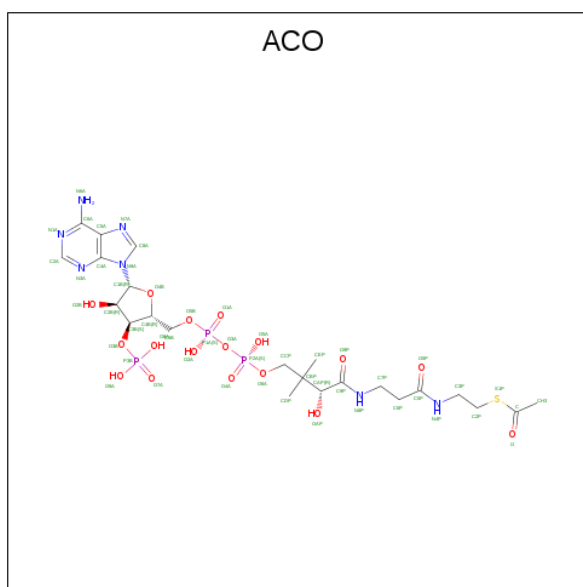
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	N	O	S				Se
1	A	326	Total	C	N	O	S	Se	0	2	0
			2464	1547	446	459	1	11			
1	B	326	Total	C	N	O	S	Se	0	2	0
			2461	1544	446	459	1	11			

- Molecule 2 is ADENOSINE-3',5'-CYCLIC-MONOPHOSPHATE (three-letter code: CMP) (formula: $C_{10}H_{12}N_5O_6P$).



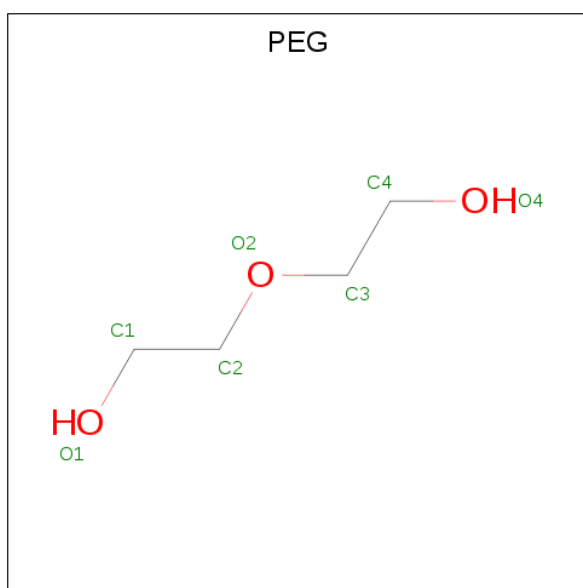
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	B	1	Total	C	N	O	P	0	0
			22	10	5	6	1		

- Molecule 3 is ACETYL COENZYME *A (three-letter code: ACO) (formula: $C_{23}H_{38}N_7O_{17}P_3S$).



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

- Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



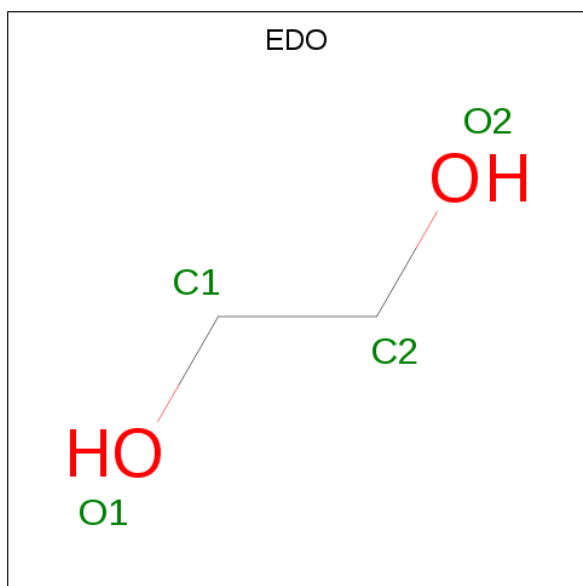
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	7	4	3	0	0
4	B	1	7	4	3	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			4	2	2		

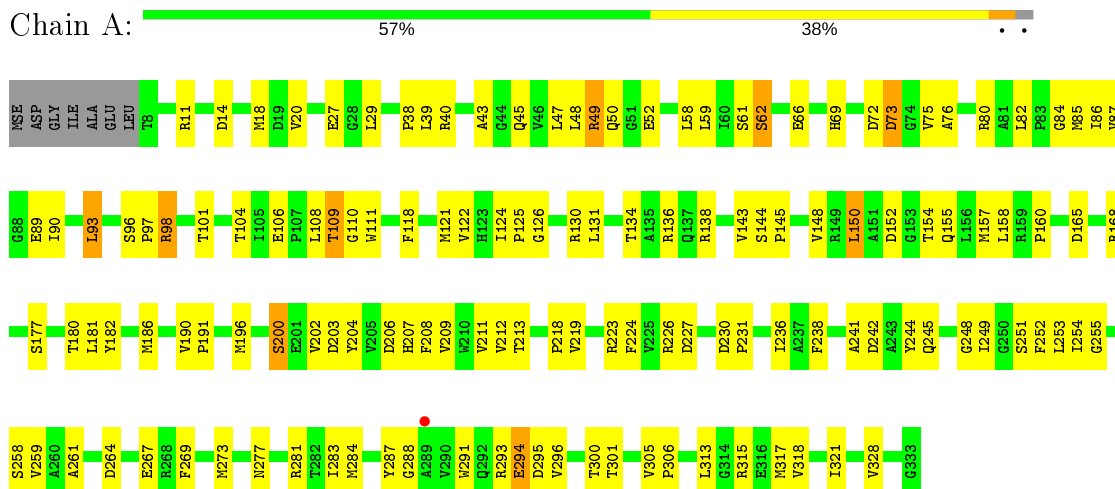
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	40	Total	O	0	0
			40	40		
6	B	43	Total	O	0	0
			43	43		

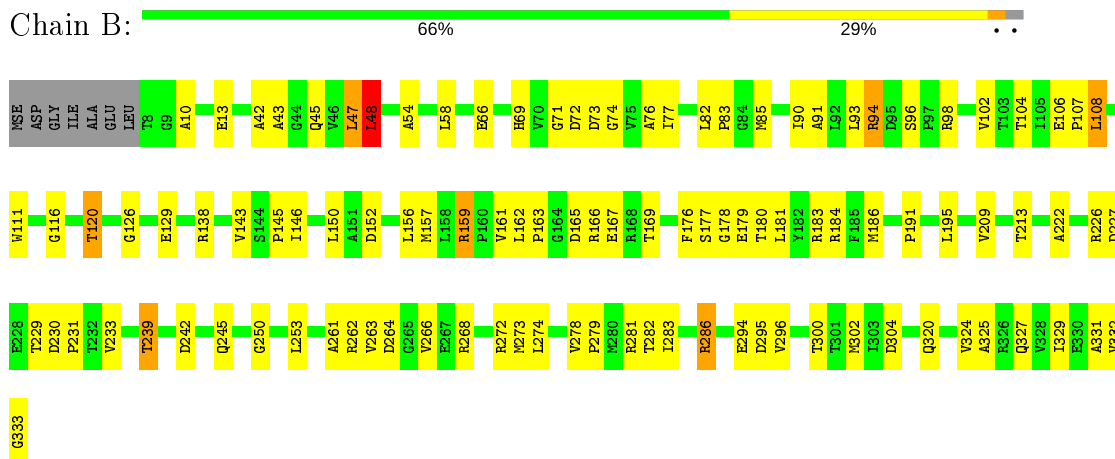
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: LYSINE ACETYLTRANSFERASE



- Molecule 1: LYSINE ACETYLTRANSFERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.66Å 50.15Å 110.10Å 90.00° 106.63° 90.00°	Depositor
Resolution (Å)	39.88 – 2.81 39.88 – 2.81	Depositor EDS
% Data completeness (in resolution range)	93.6 (39.88-2.81) 97.8 (39.88-2.81)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.81Å)	Xtrriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.199 , 0.267 0.194 , 0.266	Depositor DCC
R_{free} test set	904 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	46.7	Xtrriage
Anisotropy	0.226	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 57.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.018 for h,-k,-h-l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5179	wwPDB-VP
Average B, all atoms (Å ²)	38.0	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 31.81 % 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 1.0415e-03. 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: PEG, ACO, EDO, CMP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.35	0/2505	0.57	0/3390
1	B	0.36	0/2497	0.59	2/3376 (0.1%)
All	All	0.36	0/5002	0.58	2/6766 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	108	LEU	CA-CB-CG	5.96	129.01	115.30
1	B	48	LEU	CA-CB-CG	5.40	127.72	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	109	THR	Peptide

5.2 Too-close contacts [i](#)

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	2464	0	2474	103	0
1	B	2461	0	2472	88	0
2	A	22	0	11	5	0
2	B	22	0	11	2	0
3	A	51	0	33	6	0
3	B	51	0	33	9	0
4	A	7	0	10	5	0
4	B	14	0	20	1	0
5	B	4	0	6	1	0
6	A	40	0	0	1	0
6	B	43	0	0	1	0
All	All	5179	0	5070	197	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 19.

All (197) 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)
2:B:400:CMP:C2	2:B:400:CMP:H2	0.97	1.48
2:A:400:CMP:H2	2:A:400:CMP:C2	0.97	1.48
1:A:145:PRO:HB2	1:A:157:MSE:HE3	1.38	1.04
3:B:1334:ACO:H141	3:B:1334:ACO:N8P	1.93	0.84
1:A:252:PHE:HE2	1:A:317:MSE:HE3	1.41	0.83
1:B:294:GLU:HB2	1:B:300:THR:HG23	1.61	0.82
1:A:150:LEU:HD11	1:A:154:THR:OG1	1.81	0.80
1:A:45:GLN:HG3	6:A:2008:HOH:O	1.82	0.79
1:B:143:VAL:HG13	1:B:327:GLN:NE2	1.99	0.78
1:A:85:MSE:HE3	1:A:138:ARG:HH21	1.47	0.77
1:A:294:GLU:HB2	1:A:300:THR:HG23	1.67	0.75
1:B:43:ALA:HA	1:B:104:THR:HG22	1.69	0.75
1:B:145:PRO:HB2	1:B:157:MSE:CE	2.17	0.75
1:A:252:PHE:CE2	1:A:317:MSE:HE3	2.22	0.75
1:B:286:ARG:HH11	1:B:286:ARG:HB3	1.52	0.75
1:B:272:ARG:HG2	1:B:300:THR:HG22	1.67	0.74
1:B:227:ASP:HB3	1:B:230:ASP:O	1.87	0.74
1:B:273:MSE:HE1	1:B:281:ARG:HB2	1.69	0.74
1:A:73:ASP:OD1	1:A:75:VAL:HG22	1.89	0.72
1:B:143:VAL:HG13	1:B:327:GLN:HE22	1.58	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:VAL:HG11	1:A:321:ILE:HG22	1.75	0.69
1:A:38:PRO:HA	1:A:110:GLY:O	1.94	0.67
1:A:251:SER:O	1:A:254:ILE:HG22	1.94	0.67
3:B:1334:ACO:H141	3:B:1334:ACO:HN8	1.60	0.67
1:A:61:SER:N	1:A:109:THR:O	2.28	0.66
1:A:126:GLY:O	1:A:130:ARG:HG3	1.97	0.65
1:B:104:THR:HG23	1:B:106:GLU:O	1.97	0.65
1:B:145:PRO:HB2	1:B:157:MSE:HE2	1.79	0.64
1:A:85:MSE:HE3	1:A:138:ARG:NH2	2.12	0.64
1:A:39:LEU:HB3	1:A:110:GLY:HA2	1.80	0.63
1:B:178:GLY:HA2	1:B:181:LEU:HB2	1.79	0.63
1:B:286:ARG:NH1	1:B:286:ARG:HB3	2.12	0.63
1:B:69:HIS:O	1:B:76:ALA:HA	1.99	0.63
1:B:177:SER:HB3	1:B:242:ASP:OD2	1.99	0.63
1:B:43:ALA:HA	1:B:104:THR:CG2	2.28	0.63
1:A:209:VAL:HG22	1:A:223:ARG:HB2	1.80	0.62
1:A:131:LEU:HD23	4:A:1335:PEG:H32	1.81	0.62
1:A:177:SER:HB3	1:A:242:ASP:OD2	1.98	0.62
1:B:77:ILE:N	1:B:77:ILE:HD12	2.15	0.62
1:A:281:ARG:HG3	1:A:291:TRP:CZ2	2.34	0.62
1:B:222:ALA:HB2	1:B:253:LEU:HD13	1.82	0.62
1:A:211:VAL:CG1	1:A:218:PRO:HB3	2.30	0.61
1:A:211:VAL:HG11	1:A:218:PRO:HB3	1.81	0.61
1:A:148:VAL:HG22	1:A:317:MSE:HE1	1.83	0.61
1:B:294:GLU:HB2	1:B:300:THR:CG2	2.31	0.60
1:A:213:THR:HG22	1:A:218:PRO:HA	1.83	0.60
1:B:145:PRO:HB2	1:B:157:MSE:HE3	1.82	0.60
1:B:176:PHE:O	4:B:1336:PEG:H11	2.02	0.60
1:A:255:GLY:HA3	1:A:313:LEU:HD11	1.84	0.60
1:B:90:ILE:HG22	1:B:94:ARG:HD3	1.83	0.60
1:B:191:PRO:HB3	1:B:195:LEU:HD23	1.85	0.59
1:A:11:ARG:O	1:A:14:ASP:HB2	2.02	0.59
1:B:47:LEU:O	1:B:48:LEU:HD13	2.03	0.59
1:A:134:THR:HB	4:A:1335:PEG:H31	1.85	0.58
1:A:104:THR:HG23	1:A:106:GLU:O	2.04	0.58
1:A:18:MSE:HE2	1:A:59:LEU:CD2	2.34	0.57
1:B:286:ARG:CB	1:B:286:ARG:HH11	2.15	0.56
1:A:48:LEU:O	1:A:101:THR:HG23	2.05	0.56
1:A:267:GLU:O	1:A:305:VAL:HG23	2.05	0.56
1:A:43:ALA:HA	1:A:104:THR:HG22	1.87	0.56
1:A:49:ARG:HB2	1:A:52[A]:GLU:CD	2.26	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:THR:O	1:B:286:ARG:HG2	2.06	0.56
1:A:241:ALA:O	1:A:245:GLN:HG3	2.06	0.55
1:A:273:MSE:HE1	1:A:281:ARG:HB2	1.88	0.55
1:A:283:ILE:CG1	3:A:1334:ACO:H52A	2.37	0.55
1:A:244:TYR:HB3	1:A:249:ILE:HD12	1.88	0.55
1:A:284:MSE:O	1:A:288:GLY:O	2.24	0.54
1:B:272:ARG:HG2	1:B:300:THR:CG2	2.36	0.54
1:A:157:MSE:O	1:A:212:VAL:HA	2.08	0.53
1:B:150:LEU:HB3	1:B:152:ASP:OD1	2.09	0.53
1:B:263:VAL:HG12	1:B:329:ILE:HD12	1.91	0.53
1:A:18:MSE:HE2	1:A:59:LEU:HD23	1.89	0.53
1:B:82:LEU:HB3	1:B:83:PRO:HD2	1.91	0.53
3:A:1334:ACO:H131	3:A:1334:ACO:O9P	2.07	0.52
1:A:273:MSE:HE1	1:A:277:ASN:O	2.10	0.52
1:A:40:ARG:HD3	1:A:109:THR:HG22	1.90	0.52
1:A:254:ILE:O	1:A:258:SER:OG	2.23	0.52
1:B:138:ARG:HG2	2:B:400:CMP:C6	2.45	0.52
1:A:196:MSE:O	1:A:200:SER:HB2	2.10	0.52
1:A:90:ILE:CD1	2:A:400:CMP:H2'	2.40	0.52
1:A:50:GLN:O	1:A:98:ARG:O	2.28	0.52
1:A:138:ARG:HD3	2:A:400:CMP:C4	2.44	0.52
1:B:183:ARG:HB2	3:B:1334:ACO:H61	1.91	0.52
1:A:87:VAL:HG23	2:A:400:CMP:H1'	1.91	0.52
1:B:42:ALA:H	1:B:45[A]:GLN:NE2	2.08	0.52
1:B:159:ARG:NH2	1:B:165:ASP:OD2	2.43	0.51
1:A:58:LEU:HD12	1:A:111:TRP:O	2.10	0.51
1:A:202:VAL:HB	1:A:207:HIS:HB3	1.92	0.51
1:B:183:ARG:CB	3:B:1334:ACO:H61	2.41	0.51
1:B:166:ARG:HG3	1:B:167:GLU:N	2.25	0.51
1:A:145:PRO:HB2	1:A:157:MSE:CE	2.26	0.51
1:A:136:ARG:HD2	1:A:203:ASP:O	2.10	0.51
1:B:161:VAL:CG1	1:B:209:VAL:HB	2.41	0.51
1:A:72:ASP:OD1	1:B:262:ARG:HD2	2.11	0.51
1:A:160:PRO:HG2	1:A:328:VAL:HG21	1.92	0.50
1:B:279:PRO:O	1:B:283:ILE:HG13	2.11	0.50
1:B:263:VAL:CG1	1:B:329:ILE:HD12	2.41	0.50
1:A:39:LEU:O	1:A:110:GLY:CA	2.60	0.50
1:A:182:TYR:O	1:A:186:MSE:HA	2.12	0.50
1:B:146:ILE:HG23	1:B:320:GLN:OE1	2.11	0.50
1:A:96:SER:HB2	1:A:97:PRO:CD	2.42	0.49
1:B:184:ARG:HE	1:B:239:THR:HG22	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:GLU:H	1:B:13:GLU:CD	2.16	0.49
1:B:250:GLY:N	3:B:1334:ACO:O4A	2.45	0.49
1:A:226:ARG:HG3	1:A:231:PRO:HA	1.93	0.48
1:A:261:ALA:HB2	1:A:269:PHE:HZ	1.79	0.48
1:A:29:LEU:HD23	1:A:124:ILE:HD13	1.95	0.48
1:A:69:HIS:O	1:A:76:ALA:HA	2.13	0.48
1:B:304:ASP:OD1	5:B:1337:EDO:H12	2.14	0.48
1:B:10:ALA:HB2	1:B:111:TRP:CE3	2.48	0.48
1:B:186:MSE:HG2	1:B:186:MSE:O	2.13	0.48
1:A:86:ILE:HG23	1:A:89:GLU:OE1	2.14	0.47
1:A:18:MSE:HE2	1:A:59:LEU:HB3	1.97	0.47
1:A:118:PHE:CD1	1:A:121:MSE:HE3	2.50	0.47
1:B:91:ALA:HB1	1:B:96:SER:O	2.15	0.47
1:B:43:ALA:HB2	1:B:107:PRO:HD3	1.97	0.47
1:A:39:LEU:O	1:A:110:GLY:HA2	2.15	0.46
1:B:278:VAL:HG11	3:B:1334:ACO:H2A	1.97	0.46
1:B:163:PRO:O	1:B:166:ARG:HG2	2.15	0.46
1:B:295:ASP:OD1	1:B:296:VAL:N	2.48	0.46
1:A:39:LEU:N	1:A:110:GLY:O	2.43	0.46
1:B:116:GLY:O	1:B:120:THR:HG23	2.14	0.46
1:B:159:ARG:HH22	1:B:165:ASP:CG	2.18	0.46
1:B:156:LEU:HB3	1:B:213:THR:O	2.15	0.46
1:B:273:MSE:HE2	1:B:273:MSE:HB2	1.72	0.46
1:A:49:ARG:O	1:A:98:ARG:HB2	2.16	0.46
1:A:208:PHE:HB3	1:A:224:PHE:CE2	2.51	0.46
1:A:300:THR:O	1:A:301:THR:HB	2.16	0.46
1:B:245:GLN:HG2	3:B:1334:ACO:H143	1.98	0.45
1:B:294:GLU:HG3	1:B:300:THR:HG21	1.98	0.45
1:A:203:ASP:O	1:A:204:TYR:HB2	2.16	0.45
1:B:71:GLY:O	1:B:72:ASP:C	2.54	0.45
1:B:126:GLY:HA2	1:B:129:GLU:OE1	2.17	0.45
1:A:259:VAL:CG1	1:A:321:ILE:HG22	2.45	0.45
1:B:227:ASP:OD1	1:B:229:THR:N	2.49	0.45
1:B:261:ALA:HB1	1:B:266:VAL:HB	1.98	0.45
1:B:54:ALA:HB1	1:B:98:ARG:CZ	2.46	0.45
1:B:180:THR:HG23	3:B:1334:ACO:H72	1.98	0.45
1:B:10:ALA:HB2	1:B:111:TRP:CZ3	2.52	0.45
1:B:143:VAL:HG11	1:B:324:VAL:HG13	1.98	0.45
1:A:287:TYR:CD1	1:A:306:PRO:HG3	2.52	0.45
1:B:186:MSE:HE2	1:B:274:LEU:HD21	1.99	0.45
1:A:18:MSE:HE3	1:A:84:GLY:C	2.37	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:LEU:O	1:A:49:ARG:HG2	2.16	0.44
1:B:272:ARG:HD3	6:B:2031:HOH:O	2.17	0.44
1:B:77:ILE:N	1:B:77:ILE:CD1	2.81	0.44
1:A:165:ASP:HA	1:A:168:ARG:HB2	2.00	0.44
1:A:224:PHE:CD1	1:A:224:PHE:C	2.90	0.44
1:A:283:ILE:HG13	3:A:1334:ACO:H52A	1.99	0.44
1:A:295:ASP:OD1	1:A:296:VAL:N	2.42	0.44
1:B:66:GLU:O	1:B:102:VAL:HA	2.17	0.44
1:A:124:ILE:HA	1:A:125:PRO:HD3	1.89	0.44
3:B:1334:ACO:H131	3:B:1334:ACO:HN8	1.82	0.43
1:A:150:LEU:HD13	1:A:152:ASP:OD1	2.17	0.43
1:A:39:LEU:O	1:A:110:GLY:HA3	2.19	0.43
1:B:73:ASP:HA	1:B:74:GLY:HA2	1.65	0.43
3:A:1334:ACO:CDP	3:A:1334:ACO:O9P	2.66	0.43
1:A:253:LEU:HA	1:A:253:LEU:HD23	1.77	0.43
1:B:166:ARG:HG3	1:B:167:GLU:H	1.82	0.43
1:B:82:LEU:O	1:B:85:MSE:HB2	2.19	0.43
1:B:268:ARG:NH2	1:B:302:MSE:HE3	2.34	0.43
1:B:325:ALA:O	1:B:329:ILE:HG13	2.18	0.43
1:B:165:ASP:O	1:B:169:THR:HG23	2.18	0.43
1:A:190:VAL:HA	1:A:191:PRO:HD3	1.91	0.43
1:A:96:SER:HB2	1:A:97:PRO:HD2	2.01	0.43
1:B:45[B]:GLN:NE2	1:B:45[B]:GLN:HA	2.34	0.43
1:B:94:ARG:HG3	1:B:96:SER:OG	2.19	0.43
1:A:212:VAL:O	1:A:219:VAL:N	2.50	0.42
1:A:255:GLY:CA	1:A:313:LEU:HD11	2.49	0.42
1:A:315:ARG:O	1:A:318:VAL:HG22	2.19	0.42
1:A:227:ASP:HB3	1:A:230:ASP:O	2.20	0.42
1:B:327:GLN:O	1:B:331:ALA:HB2	2.18	0.42
1:A:85:MSE:HE1	2:A:400:CMP:H2	2.01	0.42
1:A:186:MSE:HB2	1:A:186:MSE:HE3	1.89	0.42
1:B:332:VAL:O	1:B:333[A]:GLY:C	2.58	0.42
1:A:93:LEU:HD13	1:A:118:PHE:HZ	1.84	0.41
1:A:82:LEU:H	1:A:85:MSE:SE	2.52	0.41
1:B:58:LEU:HD12	1:B:111:TRP:O	2.21	0.41
1:B:82:LEU:HB3	1:B:83:PRO:CD	2.49	0.41
1:A:62:SER:O	1:A:108:LEU:HA	2.20	0.41
1:A:86:ILE:HD13	1:A:89:GLU:HB3	2.02	0.41
1:A:248:GLY:HA2	3:A:1334:ACO:O2A	2.20	0.41
1:B:162:LEU:O	1:B:163:PRO:C	2.59	0.41
1:A:230:ASP:HA	1:A:231:PRO:HD2	1.72	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:261:ALA:HB2	1:A:269:PHE:CZ	2.54	0.41
1:B:286:ARG:HA	1:B:286:ARG:HD2	1.85	0.41
1:B:177:SER:OG	1:B:179:GLU:HB3	2.20	0.41
1:B:226:ARG:CG	1:B:231:PRO:HA	2.50	0.41
1:A:131:LEU:HA	4:A:1335:PEG:H32	2.03	0.41
1:A:144:SER:HA	1:A:145:PRO:HD3	1.89	0.41
1:A:131:LEU:HA	4:A:1335:PEG:C3	2.51	0.40
1:B:226:ARG:HG3	1:B:231:PRO:HA	2.02	0.40
1:A:203:ASP:OD1	1:A:206:ASP:N	2.52	0.40
1:A:134:THR:CB	4:A:1335:PEG:H31	2.51	0.40
1:A:180:THR:HG23	3:A:1334:ACO:H72	2.04	0.40
1:A:236:ILE:HD12	1:A:238:PHE:CZ	2.56	0.40
1:B:253:LEU:HA	1:B:253:LEU:HD23	1.76	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	326/333 (98%)	312 (96%)	14 (4%)	0	100	100
1	B	325/333 (98%)	316 (97%)	9 (3%)	0	100	100
All	All	651/666 (98%)	628 (96%)	23 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	258/249 (104%)	238 (92%)	20 (8%)	12	35
1	B	257/249 (103%)	246 (96%)	11 (4%)	29	62
All	All	515/498 (103%)	484 (94%)	31 (6%)	18	48

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	VAL
1	A	27	GLU
1	A	47	LEU
1	A	49	ARG
1	A	62	SER
1	A	66	GLU
1	A	73	ASP
1	A	80	ARG
1	A	93	LEU
1	A	98	ARG
1	A	122	VAL
1	A	143	VAL
1	A	150	LEU
1	A	155	GLN
1	A	158	LEU
1	A	181	LEU
1	A	200	SER
1	A	264	ASP
1	A	293	ARG
1	A	294	GLU
1	B	47	LEU
1	B	48	LEU
1	B	93	LEU
1	B	94	ARG
1	B	108	LEU
1	B	120	THR
1	B	159	ARG
1	B	233	VAL
1	B	239	THR
1	B	264	ASP
1	B	286	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no

such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

8 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	CMP	A	400	-	22,25,25	1.55	5 (22%)	24,39,39	1.78	5 (20%)
2	CMP	B	400	-	22,25,25	1.53	5 (22%)	24,39,39	1.60	5 (20%)
3	ACO	B	1334	-	45,53,53	1.68	9 (20%)	56,79,79	1.56	12 (21%)
4	PEG	B	1336	-	6,6,6	0.53	0	5,5,5	1.76	2 (40%)
3	ACO	A	1334	-	45,53,53	1.72	10 (22%)	56,79,79	1.64	9 (16%)
4	PEG	B	1335	-	6,6,6	0.61	0	5,5,5	1.50	1 (20%)
4	PEG	A	1335	-	6,6,6	0.59	0	5,5,5	1.59	0
5	EDO	B	1337	-	3,3,3	0.55	0	2,2,2	0.13	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	CMP	A	400	-	-	0/0/31/31	0/4/4/4
2	CMP	B	400	-	-	0/0/31/31	0/4/4/4
3	ACO	B	1334	-	1/1/12/14	17/47/67/67	0/3/3/3
4	PEG	B	1336	-	-	2/4/4/4	-
3	ACO	A	1334	-	1/1/12/14	23/47/67/67	0/3/3/3
4	PEG	B	1335	-	-	2/4/4/4	-
4	PEG	A	1335	-	-	3/4/4/4	-
5	EDO	B	1337	-	-	1/1/1/1	-

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1334	ACO	C5P-N4P	5.88	1.46	1.33
3	B	1334	ACO	C5P-N4P	5.72	1.46	1.33
3	A	1334	ACO	C9P-N8P	4.68	1.43	1.33
3	B	1334	ACO	C9P-N8P	4.41	1.43	1.33
2	A	400	CMP	P-O5'	3.47	1.61	1.57
2	A	400	CMP	P-O3'	3.42	1.63	1.57
2	B	400	CMP	P-O5'	3.21	1.61	1.57
2	B	400	CMP	P-O3'	3.07	1.62	1.57
3	A	1334	ACO	C6A-N6A	3.03	1.45	1.34
3	A	1334	ACO	C2B-C1B	-3.00	1.49	1.53
3	B	1334	ACO	C6A-N6A	3.00	1.45	1.34
3	B	1334	ACO	O2B-C2B	-2.81	1.36	1.43
3	B	1334	ACO	C5B-C4B	-2.78	1.42	1.51
3	A	1334	ACO	C2B-C3B	-2.75	1.46	1.52
3	A	1334	ACO	O2B-C2B	-2.74	1.36	1.43
3	B	1334	ACO	C2B-C1B	-2.61	1.49	1.53
2	A	400	CMP	O3'-C3'	-2.57	1.40	1.44
3	A	1334	ACO	P3B-O3B	-2.55	1.54	1.59
3	B	1334	ACO	C2B-C3B	-2.48	1.47	1.52
3	A	1334	ACO	C5B-C4B	-2.46	1.43	1.51
2	B	400	CMP	C5-C4	2.45	1.47	1.40
2	A	400	CMP	O5'-C5'	-2.34	1.42	1.46
2	B	400	CMP	O3'-C3'	-2.28	1.41	1.44
2	B	400	CMP	O5'-C5'	-2.26	1.43	1.46
3	A	1334	ACO	O3B-C3B	-2.21	1.36	1.44
3	A	1334	ACO	OAP-CAP	-2.17	1.38	1.42
3	B	1334	ACO	P3B-O3B	-2.17	1.55	1.59
2	A	400	CMP	C5-C4	2.14	1.46	1.40
3	B	1334	ACO	O3B-C3B	-2.04	1.36	1.44

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1334	ACO	O6A-CCP-CBP	5.87	119.98	110.55
3	A	1334	ACO	N3A-C2A-N1A	-4.73	121.28	128.68
2	A	400	CMP	O5'-P-O3'	4.47	111.84	105.68
3	B	1334	ACO	N3A-C2A-N1A	-4.32	121.92	128.68
2	A	400	CMP	O3'-P-O1P	-3.73	102.40	110.39
3	B	1334	ACO	C6P-C5P-N4P	3.64	122.54	116.42
2	B	400	CMP	O3'-P-O1P	-3.52	102.85	110.39
3	B	1334	ACO	P2A-O3A-P1A	-3.49	120.86	132.83
3	A	1334	ACO	C7P-C6P-C5P	3.41	118.04	112.36
3	B	1334	ACO	O5P-C5P-N4P	-3.37	116.64	123.01
3	B	1334	ACO	C3P-N4P-C5P	-3.33	116.66	122.84
2	B	400	CMP	O2P-P-O1P	3.26	118.94	108.73
2	A	400	CMP	N3-C2-N1	-3.21	123.67	128.68
2	A	400	CMP	O2P-P-O1P	3.18	118.68	108.73
3	A	1334	ACO	CDP-CBP-CAP	3.15	114.27	108.82
2	B	400	CMP	N3-C2-N1	-3.10	123.84	128.68
3	A	1334	ACO	O5B-C5B-C4B	3.02	119.40	108.99
3	B	1334	ACO	C1B-N9A-C4A	-2.88	121.58	126.64
3	B	1334	ACO	C7P-C6P-C5P	2.77	116.97	112.36
3	A	1334	ACO	C7P-N8P-C9P	-2.73	117.72	122.59
3	B	1334	ACO	C6P-C7P-N8P	-2.52	106.82	111.90
2	B	400	CMP	O5'-P-O3'	2.52	109.14	105.68
3	B	1334	ACO	C7P-N8P-C9P	-2.51	118.11	122.59
2	A	400	CMP	C4-C5-N7	-2.35	106.95	109.40
3	A	1334	ACO	C3P-N4P-C5P	-2.33	118.51	122.84
3	B	1334	ACO	CEP-CBP-CDP	2.33	113.91	109.17
4	B	1336	PEG	O2-C3-C4	2.28	120.07	110.07
2	B	400	CMP	C4-C5-N7	-2.26	107.04	109.40
3	B	1334	ACO	O5B-C5B-C4B	2.22	116.62	108.99
3	A	1334	ACO	P2A-O3A-P1A	-2.19	125.32	132.83
4	B	1336	PEG	O2-C2-C1	2.16	119.58	110.07
3	B	1334	ACO	C3B-C2B-C1B	2.11	104.57	99.89
3	A	1334	ACO	OAP-CAP-CBP	2.11	115.21	110.25
4	B	1335	PEG	O2-C2-C1	2.01	118.92	110.07

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	1334	ACO	CAP
3	A	1334	ACO	CAP

All (48) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1334	ACO	O4B-C4B-C5B-O5B
3	B	1334	ACO	C5B-O5B-P1A-O3A
3	B	1334	ACO	C9P-CAP-CBP-CCP
3	B	1334	ACO	O9P-C9P-CAP-CBP
3	B	1334	ACO	N8P-C9P-CAP-CBP
3	B	1334	ACO	C5P-C6P-C7P-N8P
3	A	1334	ACO	C3B-O3B-P3B-O7A
3	A	1334	ACO	C5B-O5B-P1A-O3A
3	A	1334	ACO	P1A-O3A-P2A-O6A
3	A	1334	ACO	CDP-CBP-CCP-O6A
3	A	1334	ACO	CEP-CBP-CCP-O6A
3	A	1334	ACO	CAP-CBP-CCP-O6A
3	A	1334	ACO	OAP-CAP-CBP-CCP
3	A	1334	ACO	OAP-CAP-CBP-CDP
3	A	1334	ACO	O9P-C9P-CAP-OAP
3	A	1334	ACO	C5P-C6P-C7P-N8P
3	B	1334	ACO	C3B-C4B-C5B-O5B
4	B	1336	PEG	O1-C1-C2-O2
4	B	1336	PEG	O2-C3-C4-O4
4	B	1335	PEG	O2-C3-C4-O4
4	A	1335	PEG	O2-C3-C4-O4
3	B	1334	ACO	O9P-C9P-CAP-OAP
3	B	1334	ACO	OAP-CAP-CBP-CDP
3	A	1334	ACO	OAP-CAP-CBP-CEP
3	A	1334	ACO	N8P-C9P-CAP-OAP
4	B	1335	PEG	C4-C3-O2-C2
4	A	1335	PEG	C4-C3-O2-C2
3	B	1334	ACO	C9P-CAP-CBP-CDP
3	B	1334	ACO	C3B-O3B-P3B-O9A
3	B	1334	ACO	C5B-O5B-P1A-O1A
3	B	1334	ACO	C5B-O5B-P1A-O2A
3	A	1334	ACO	C5B-O5B-P1A-O2A
3	A	1334	ACO	CCP-O6A-P2A-O4A
4	A	1335	PEG	C1-C2-O2-C3
3	B	1334	ACO	OAP-CAP-CBP-CCP
3	A	1334	ACO	O9P-C9P-CAP-CBP
3	B	1334	ACO	CEP-CBP-CCP-O6A
3	A	1334	ACO	N8P-C9P-CAP-CBP
3	A	1334	ACO	C3P-C2P-S1P-C
5	B	1337	EDO	O1-C1-C2-O2
3	A	1334	ACO	O4B-C4B-C5B-O5B
3	B	1334	ACO	C9P-CAP-CBP-CEP
3	A	1334	ACO	C9P-CAP-CBP-CDP

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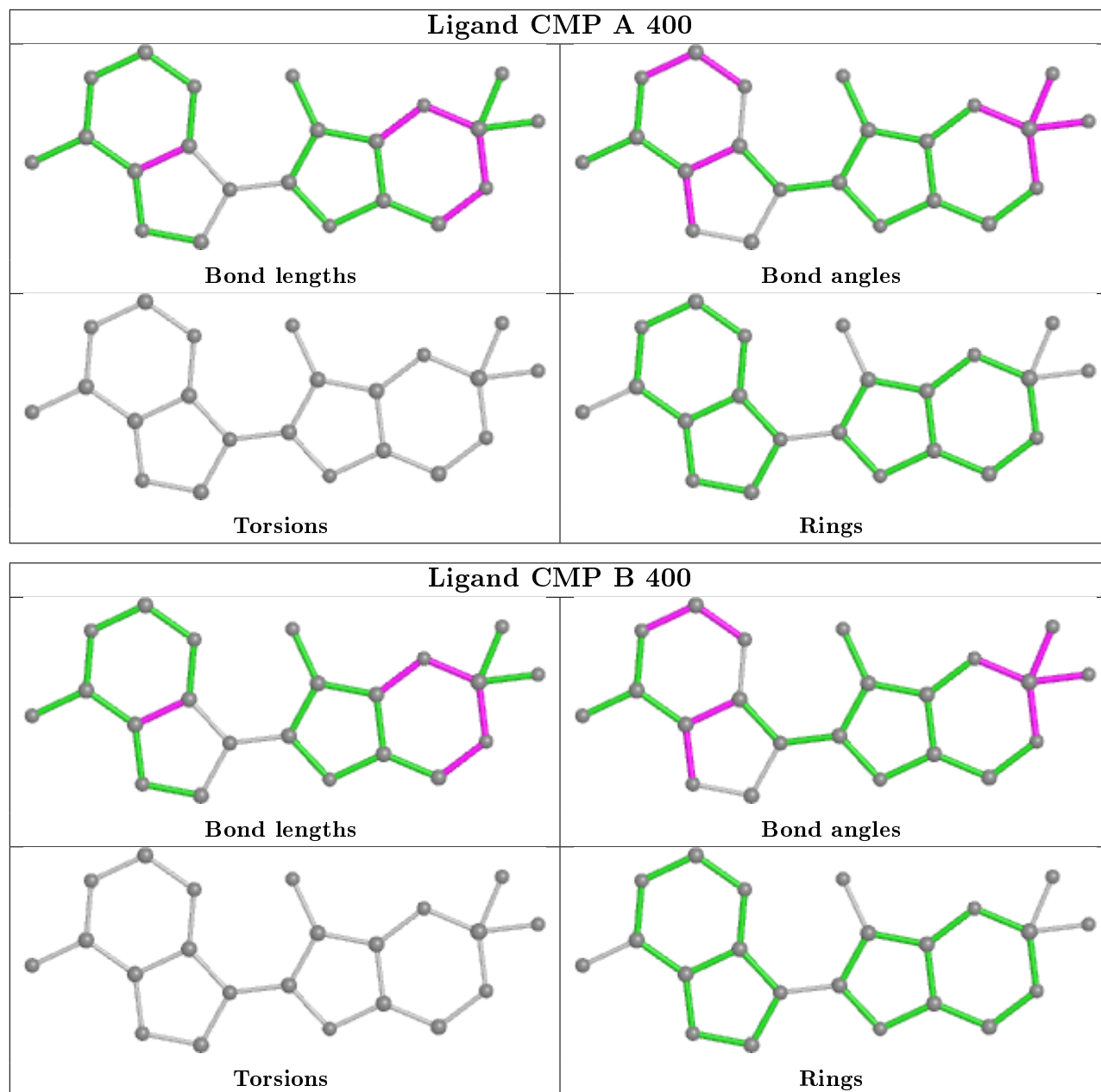
Mol	Chain	Res	Type	Atoms
3	A	1334	ACO	C9P-CAP-CBP-CEP
3	A	1334	ACO	CCP-O6A-P2A-O3A
3	A	1334	ACO	P2A-O3A-P1A-O2A
3	A	1334	ACO	CBP-CCP-O6A-P2A
3	B	1334	ACO	CH3-C-S1P-C2P

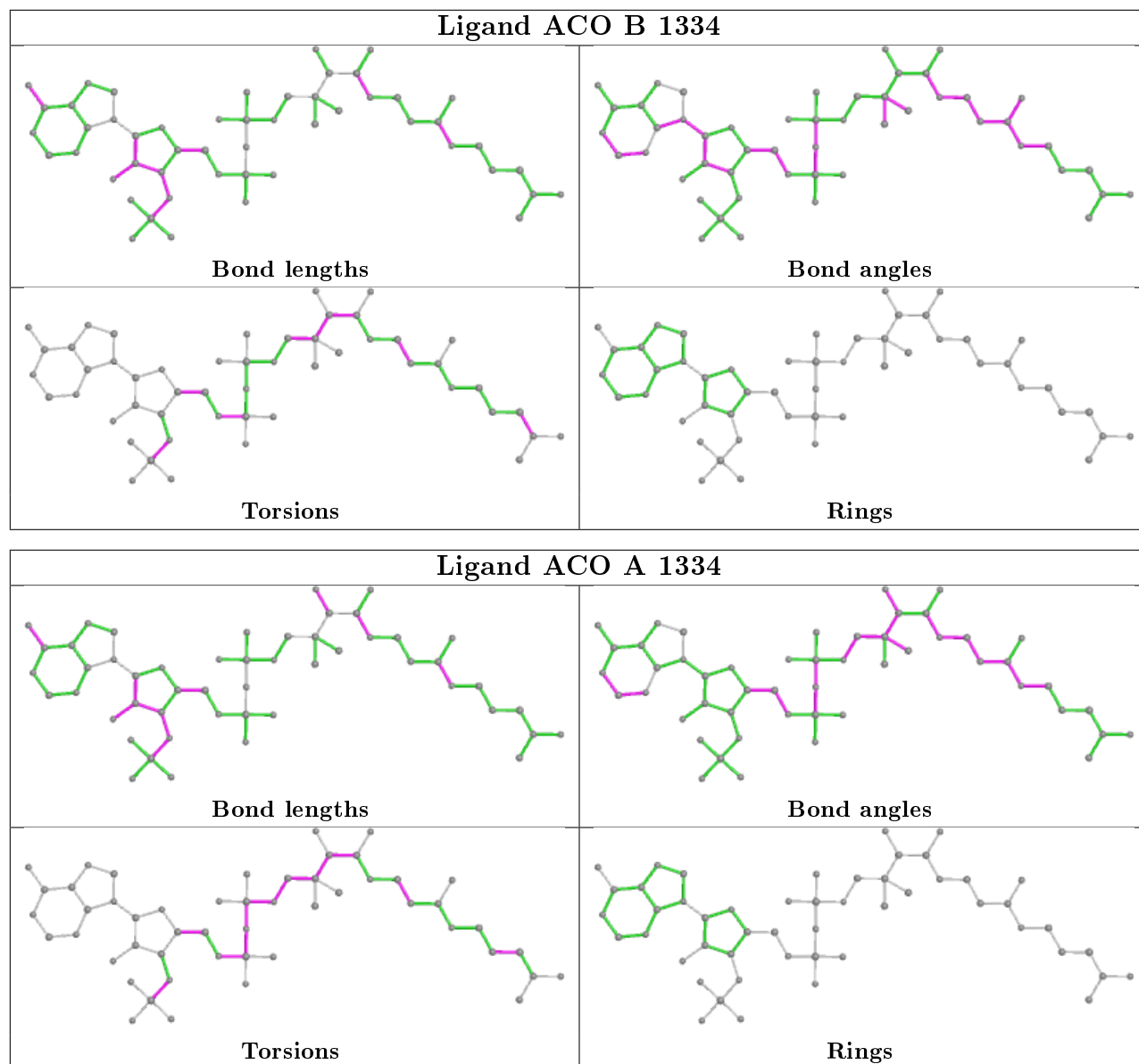
There are no ring outliers.

7 monomers are involved in 29 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	400	CMP	5	0
2	B	400	CMP	2	0
3	B	1334	ACO	9	0
4	B	1336	PEG	1	0
3	A	1334	ACO	6	0
4	A	1335	PEG	5	0
5	B	1337	EDO	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	315/333 (94%)	-0.32	1 (0%) 94 93	18, 38, 62, 80	0
1	B	315/333 (94%)	-0.38	0 100 100	17, 34, 60, 82	0
All	All	630/666 (94%)	-0.35	1 (0%) 95 94	17, 36, 62, 82	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	289	ALA	2.4

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 carbohydrates 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.

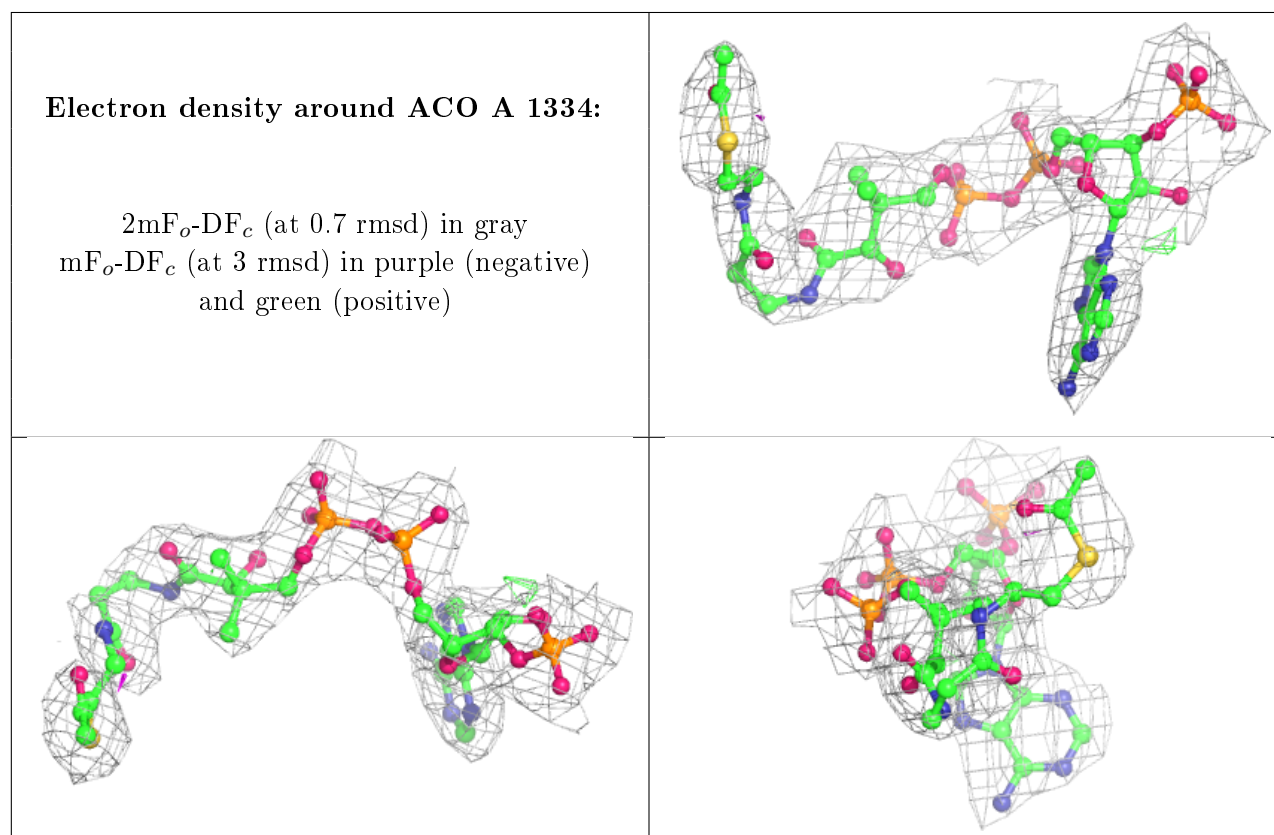
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	PEG	B	1336	7/7	0.86	0.24	36,38,43,44	0
4	PEG	A	1335	7/7	0.87	0.43	49,52,61,61	0
4	PEG	B	1335	7/7	0.91	0.28	49,52,53,53	0
3	ACO	A	1334	51/51	0.94	0.18	21,52,72,73	0

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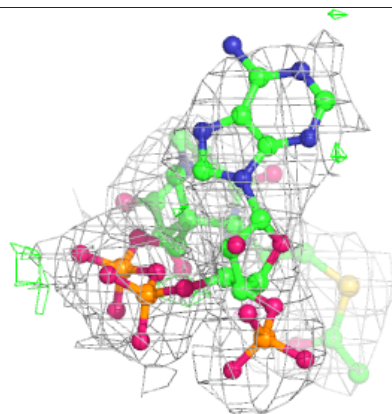
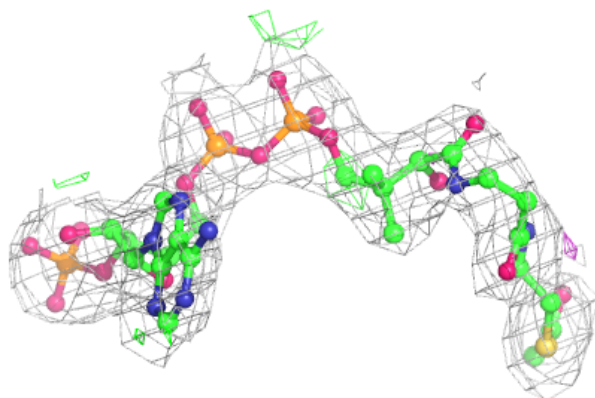
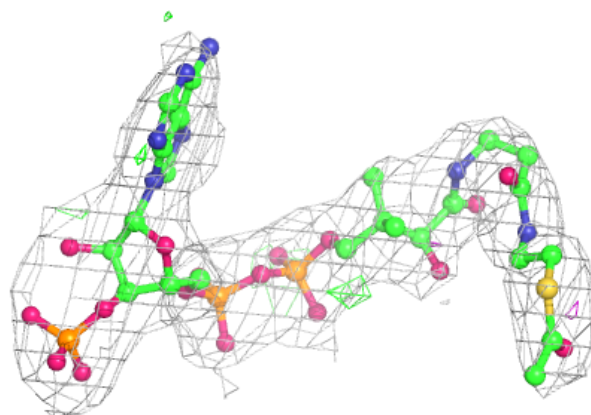
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	ACO	B	1334	51/51	0.95	0.17	26,39,70,74	0
5	EDO	B	1337	4/4	0.95	0.22	37,39,39,40	0
2	CMP	A	400	22/22	0.98	0.15	20,25,28,31	0
2	CMP	B	400	22/22	0.98	0.14	29,32,35,37	0

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.

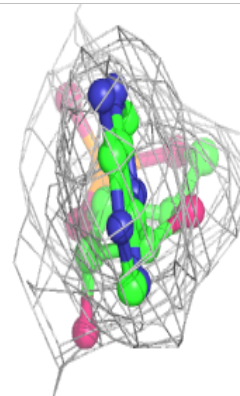
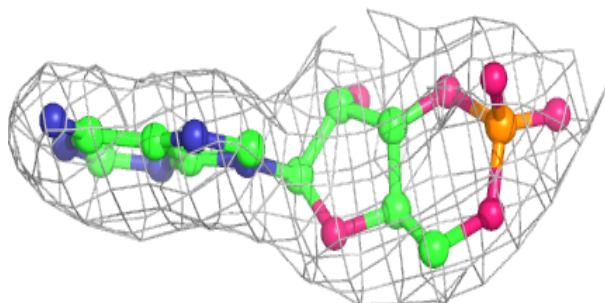
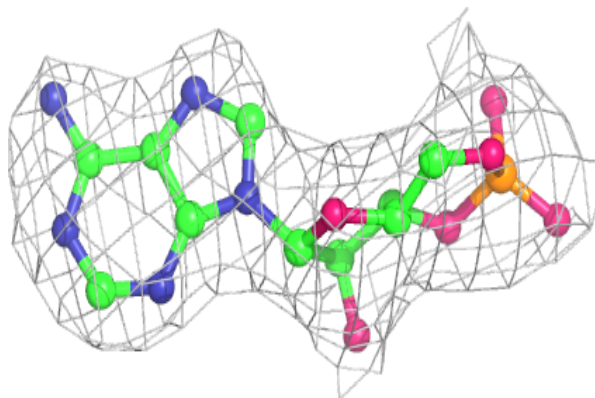


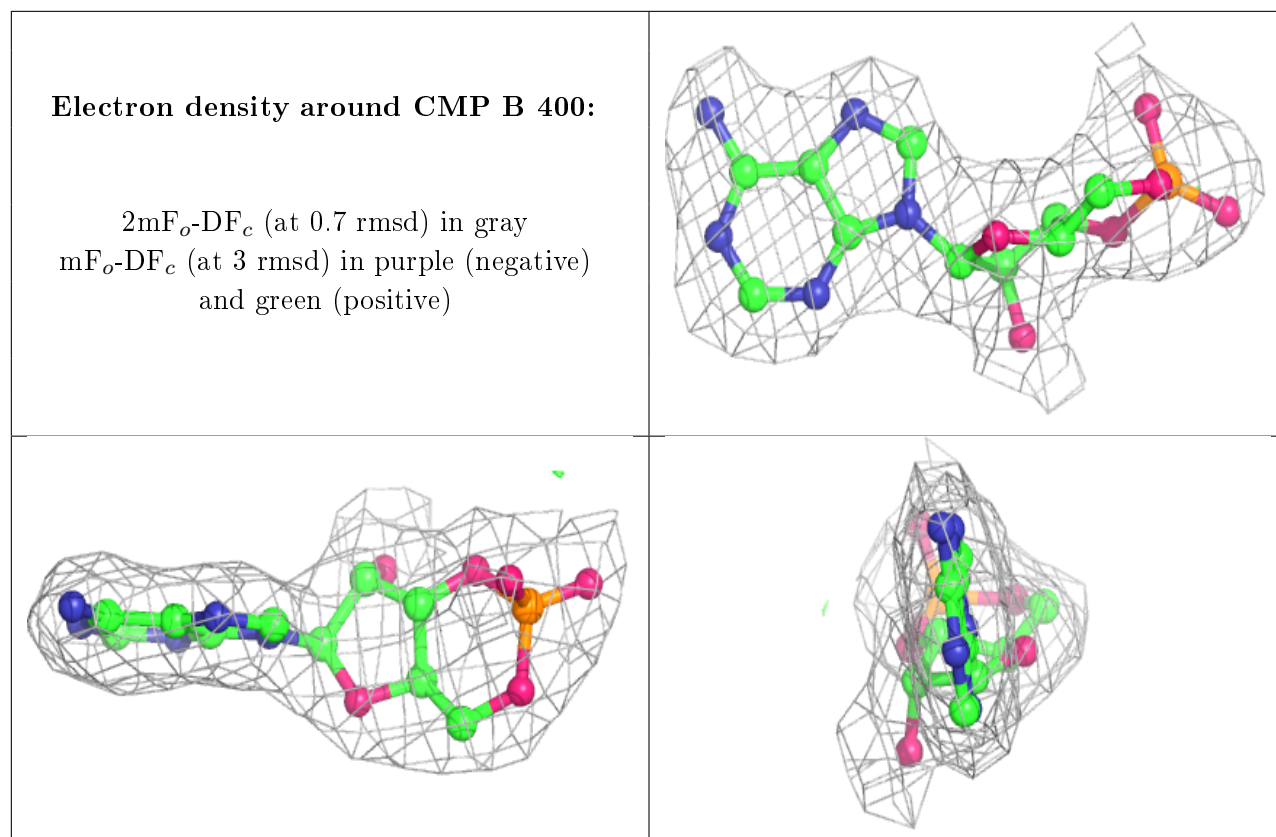
Electron density around ACO B 1334:

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

**Electron density around CMP A 400:**

$2mF_o-DF_c$ (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.