Full wwPDB X-ray Structure Validation Report (i)

Aug 28, 2023 – 12:12 AM EDT

PDB ID	:	3KLO
Title	:	Vibrio cholerae VpsT bound to c-di-GMP
Authors	:	Krasteva, P.V.; Navarro, V.A.S.; Sondermann, H.
Deposited on	:	2009-11-08
Resolution	:	2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (i)) were used in the production of this report:

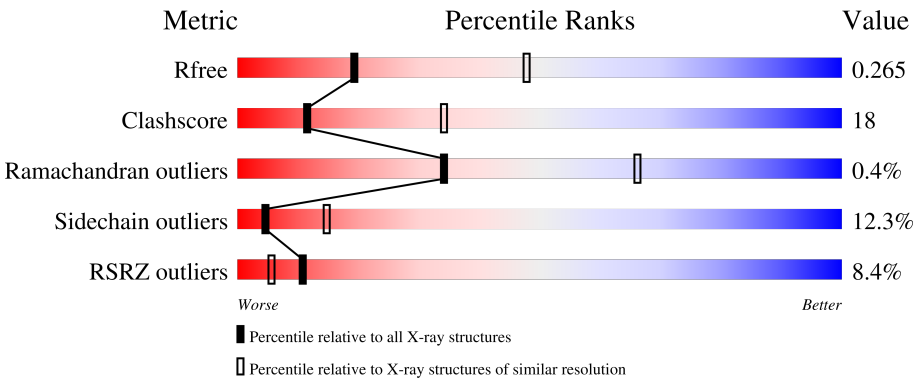
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	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 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 <=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	225	62% 28% 7% •
1	B	225	4% 62% 30% • •
1	C	225	12% 68% 20% 7% 5%
1	D	225	14% 51% 21% 5% 23%

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	TAR	B	226	-	-	X	X
3	TAR	D	226	X	-	-	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6902 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.

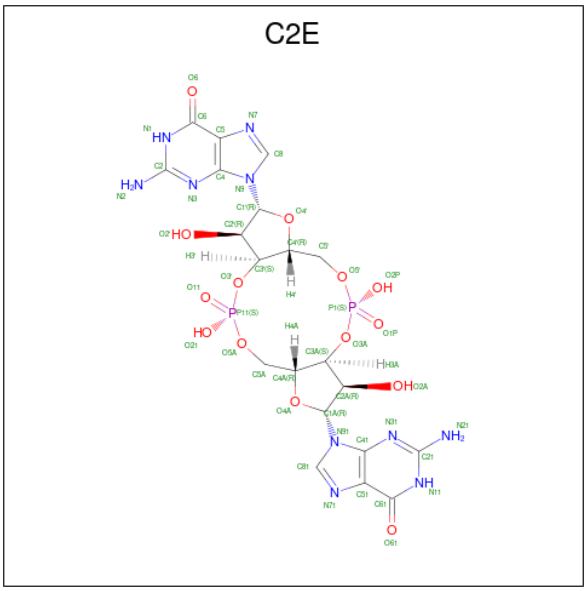
- Molecule 1 is a protein called Transcriptional regulator VpsT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	217	1756	1119	298	329	10	0	0	0
1	B	216	1748	1115	297	326	10	0	0	0
1	C	213	1728	1101	294	323	10	0	0	0
1	D	173	1417	905	241	262	9	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

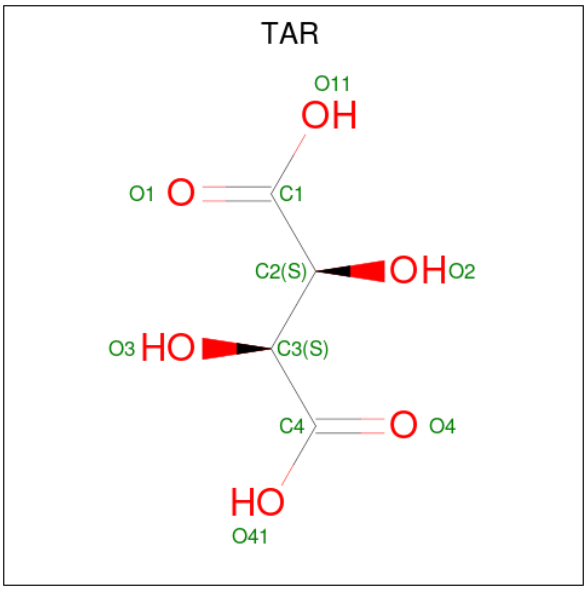
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q9KKZ8
B	0	SER	-	expression tag	UNP Q9KKZ8
C	0	SER	-	expression tag	UNP Q9KKZ8
D	0	SER	-	expression tag	UNP Q9KKZ8

- Molecule 2 is 9,9'-[(2R,3R,3aS,5S,7aR,9R,10R,10aS,12S,14aR)-3,5,10,12-tetrahydroxy-5,12-dioxidooctahydro-2H,7H-difuro[3,2-d:3',2'-j][1,3,7,9,2,8]tetraoxadiphosphacyclododecine-2,9-diyl]bis(2-amino-1,9-dihydro-6H-purin-6-one) (three-letter code: C2E) (formula: $C_{20}H_{24}N_{10}O_{14}P_2$).



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

- Molecule 3 is D(-)-TARTARIC ACID (three-letter code: TAR) (formula: $C_4H_6O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	B	1	10	4	6	0	0
			Total	C	O		
3	D	1	10	4	6	0	0

- Molecule 4 is water.

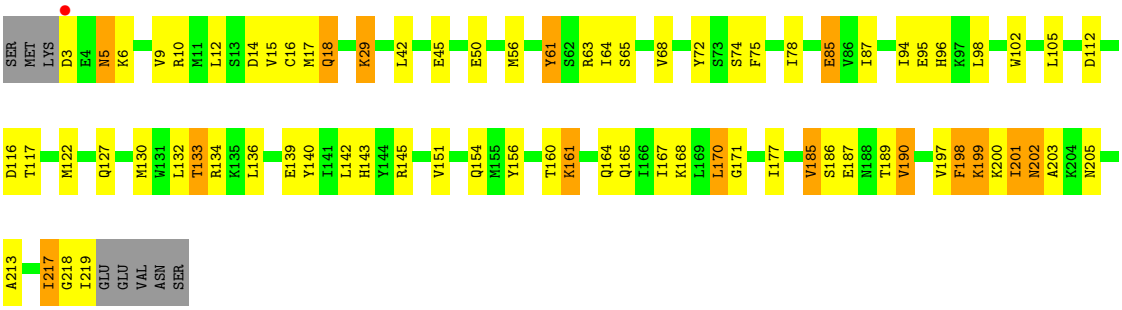
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	12	12	12	0	0
			Total	O		
4	B	19	19	19	0	0
			Total	O		
4	C	7	7	7	0	0
			Total	O		
4	D	11	11	11	0	0

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.

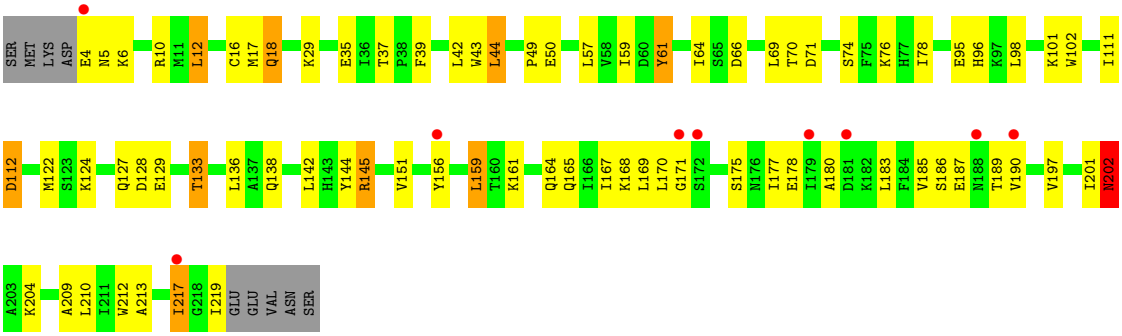
• Molecule 1: Transcriptional regulator VpsT

Chain A: 62% 28% 7%



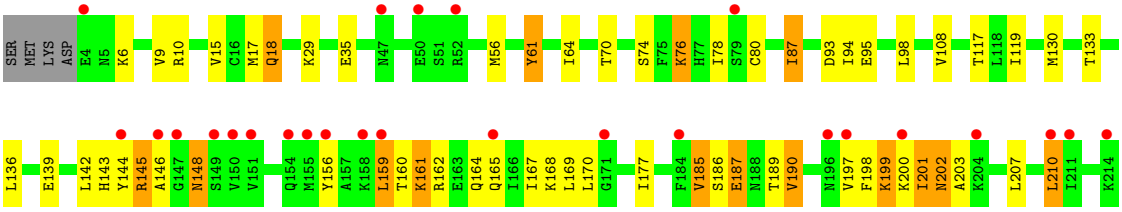
• Molecule 1: Transcriptional regulator VpsT

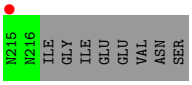
Chain B: 4% 62% 30%



• Molecule 1: Transcriptional regulator VpsT

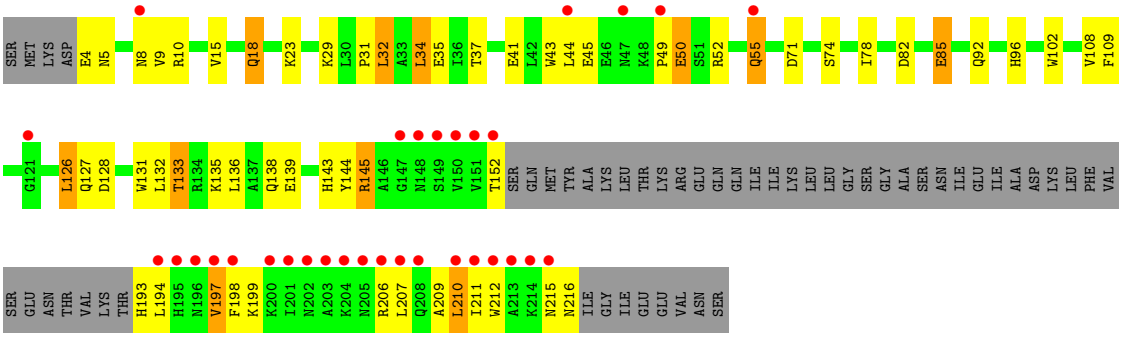
Chain C: 12% 68% 20% 7% 5%





• Molecule 1: Transcriptional regulator VpsT

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	121.70Å 121.70Å 208.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.86 – 2.80 47.86 – 2.80	Depositor EDS
% Data completeness (in resolution range)	82.8 (47.86-2.80) 98.0 (47.86-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.97 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 2009_02_15_2320_3)	Depositor
R, R_{free}	0.246 , 0.290 0.266 , 0.265	Depositor DCC
R_{free} test set	1922 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	50.1	Xtriage
Anisotropy	0.186	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 46.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	6902	wwPDB-VP
Average B, all atoms (Å^2)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 74.77 % 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 1.4350e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: C2E, TAR

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	0.41	0/1785	0.54	0/2407
1	B	0.38	0/1777	0.54	0/2396
1	C	0.31	0/1757	0.48	0/2369
1	D	0.34	0/1443	0.52	0/1947
All	All	0.36	0/6762	0.52	0/9119

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	1756	0	1796	72	0
1	B	1748	0	1792	72	0
1	C	1728	0	1767	56	0
1	D	1417	0	1435	49	0
2	A	46	0	22	0	0
2	B	46	0	22	2	0
2	C	46	0	22	1	0
2	D	46	0	22	1	0
3	B	10	0	6	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	10	0	6	0	0
4	A	12	0	0	0	0
4	B	19	0	0	2	0
4	C	7	0	0	1	0
4	D	11	0	0	0	0
All	All	6902	0	6890	242	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 18.

All (242) 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:D:145:ARG:HH11	1:D:145:ARG:HG2	1.13	1.09
1:B:145:ARG:HG2	1:B:145:ARG:HH11	1.13	1.04
1:A:18:GLN:H	1:A:18:GLN:HE21	1.05	1.01
1:B:133:THR:HG22	1:B:136:LEU:H	1.27	1.00
1:D:18:GLN:H	1:D:18:GLN:HE21	1.06	0.94
1:B:18:GLN:H	1:B:18:GLN:HE21	0.95	0.93
1:D:210:LEU:H	1:D:210:LEU:HD23	1.41	0.86
1:B:18:GLN:H	1:B:18:GLN:NE2	1.76	0.82
1:A:167:ILE:HD13	1:A:197:VAL:HG11	1.63	0.81
1:C:156:TYR:OH	1:C:168:LYS:HE2	1.81	0.81
1:B:133:THR:CG2	1:B:136:LEU:H	1.94	0.80
1:C:18:GLN:H	1:C:18:GLN:HE21	1.28	0.79
1:A:186:SER:HB3	1:A:189:THR:OG1	1.83	0.79
1:D:145:ARG:HH11	1:D:145:ARG:CG	1.94	0.78
1:D:18:GLN:H	1:D:18:GLN:NE2	1.81	0.77
1:D:4:GLU:HG3	1:D:5:ASN:H	1.48	0.77
1:B:145:ARG:HG2	1:B:145:ARG:NH1	1.94	0.77
1:B:213:ALA:O	1:B:217:ILE:HG22	1.84	0.76
1:C:145:ARG:HG2	1:C:145:ARG:HH11	1.49	0.76
1:D:8:ASN:HB3	1:D:55:GLN:HG3	1.67	0.76
1:B:177:ILE:HG13	1:B:187:GLU:HG3	1.66	0.75
1:A:133:THR:HG22	1:A:136:LEU:H	1.50	0.75
1:C:145:ARG:HH11	1:C:145:ARG:CG	1.99	0.74
1:A:133:THR:CG2	1:A:136:LEU:H	2.01	0.74
1:B:197:VAL:HG12	1:B:201:ILE:HD12	1.70	0.73
1:D:145:ARG:HG2	1:D:145:ARG:NH1	1.94	0.73
1:A:130:MET:HG3	1:B:138:GLN:HG3	1.69	0.72
1:B:4:GLU:HG2	1:B:5:ASN:H	1.54	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:HIS:CD2	1:A:139:GLU:HG3	2.25	0.72
1:C:133:THR:HG22	1:C:136:LEU:HB3	1.71	0.71
1:D:207:LEU:HA	1:D:210:LEU:HD21	1.71	0.70
2:C:225:C2E:H5'2	2:C:225:C2E:H512	1.74	0.69
1:B:4:GLU:HG2	1:B:5:ASN:N	2.08	0.69
1:C:61:TYR:O	1:C:64:ILE:HG22	1.93	0.69
1:C:9:VAL:HG22	1:C:56:MET:HB3	1.74	0.68
1:D:133:THR:HG22	1:D:136:LEU:H	1.58	0.68
1:C:177:ILE:HG13	1:C:187:GLU:HG3	1.74	0.68
1:A:127:GLN:HE22	1:C:6:LYS:HE3	1.59	0.67
1:B:43:TRP:CE2	1:B:44:LEU:HD22	2.30	0.67
1:D:31:PRO:HD2	1:D:126:LEU:HD23	1.76	0.66
1:B:133:THR:HG22	1:B:136:LEU:N	2.08	0.66
1:B:201:ILE:O	1:B:202:ASN:HB2	1.94	0.66
1:A:177:ILE:HG13	1:A:187:GLU:HG3	1.78	0.66
1:A:127:GLN:NE2	1:C:6:LYS:HE3	2.11	0.65
1:B:124:LYS:HD3	1:B:129:GLU:OE2	1.96	0.65
1:A:95:GLU:HB3	1:A:98:LEU:HG	1.78	0.65
1:C:159:LEU:HD12	1:C:164:GLN:HG2	1.78	0.65
1:A:5:ASN:C	1:A:5:ASN:HD22	1.99	0.65
1:A:18:GLN:H	1:A:18:GLN:NE2	1.86	0.65
1:B:112:ASP:N	1:B:112:ASP:OD1	2.30	0.64
1:B:161:LYS:O	1:B:165:GLN:HG3	1.96	0.64
1:C:207:LEU:HA	1:C:210:LEU:HD23	1.78	0.64
1:D:194:LEU:O	1:D:197:VAL:HG23	1.99	0.63
1:A:156:TYR:OH	1:A:168:LYS:HE3	1.98	0.63
1:B:145:ARG:HH11	1:B:145:ARG:CG	2.00	0.63
1:C:185:VAL:HG13	1:C:189:THR:HB	1.79	0.62
1:C:133:THR:HG22	1:C:136:LEU:CB	2.29	0.62
1:C:167:ILE:HD13	1:C:197:VAL:HG11	1.80	0.62
1:D:18:GLN:HE21	1:D:18:GLN:N	1.89	0.62
1:A:9:VAL:HG22	1:A:56:MET:HB3	1.80	0.61
1:C:145:ARG:HG2	1:C:145:ARG:NH1	2.14	0.61
1:C:74:SER:O	1:C:78:ILE:HG12	2.00	0.61
1:C:159:LEU:CD1	1:C:164:GLN:HG2	2.30	0.61
1:D:210:LEU:H	1:D:210:LEU:CD2	2.13	0.60
1:B:197:VAL:CG1	1:B:201:ILE:HD12	2.32	0.60
1:A:145:ARG:HA	1:B:144:TYR:CD1	2.35	0.60
1:A:134:ARG:HG3	3:B:226:TAR:O1	2.01	0.60
1:C:133:THR:CG2	1:C:136:LEU:H	2.14	0.60
1:C:161:LYS:N	1:C:161:LYS:HD2	2.17	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:186:SER:HB3	1:C:189:THR:OG1	2.02	0.59
1:B:204:LYS:HE3	4:B:245:HOH:O	2.02	0.59
1:C:95:GLU:HB3	1:C:98:LEU:HG	1.82	0.59
1:D:85:GLU:OE1	1:D:102:TRP:HE3	1.86	0.59
1:C:143:HIS:CE1	1:C:148:ASN:HB2	2.37	0.59
1:C:201:ILE:HG22	1:C:203:ALA:H	1.68	0.59
1:A:74:SER:O	1:A:78:ILE:HG12	2.03	0.59
1:C:76:LYS:HE3	1:C:80:CYS:O	2.02	0.58
1:B:124:LYS:HE3	3:B:226:TAR:H2	1.86	0.58
1:B:49:PRO:HD2	1:B:50:GLU:OE1	2.04	0.58
1:A:161:LYS:O	1:A:165:GLN:HG3	2.04	0.58
1:A:18:GLN:HE21	1:A:18:GLN:N	1.89	0.58
1:C:143:HIS:O	1:C:146:ALA:HB3	2.04	0.57
1:A:29:LYS:HG3	1:C:119:ILE:HG22	1.86	0.57
1:B:18:GLN:HE21	1:B:18:GLN:N	1.81	0.57
1:B:61:TYR:O	1:B:64:ILE:HG22	2.04	0.57
1:D:133:THR:HG22	1:D:135:LYS:N	2.21	0.56
1:A:197:VAL:C	1:A:199:LYS:H	2.09	0.55
1:C:145:ARG:HH11	1:C:145:ARG:HB3	1.71	0.55
1:B:66:ASP:HA	1:B:69:LEU:HD12	1.88	0.55
1:C:201:ILE:HG22	1:C:203:ALA:N	2.22	0.55
1:B:186:SER:HB3	1:B:189:THR:OG1	2.06	0.55
1:D:4:GLU:HG3	1:D:5:ASN:N	2.20	0.55
1:C:160:THR:O	1:C:164:GLN:HG3	2.07	0.55
1:B:59:ILE:HD13	1:B:102:TRP:CZ3	2.43	0.54
1:A:167:ILE:CG2	1:A:168:LYS:N	2.70	0.54
1:B:219:ILE:HG23	1:B:219:ILE:O	2.08	0.54
1:C:145:ARG:HH11	1:C:145:ARG:CB	2.20	0.54
1:A:133:THR:HG22	1:A:136:LEU:CB	2.38	0.54
1:A:87:ILE:HD13	1:A:105:LEU:CD1	2.38	0.53
1:C:70:THR:HG23	4:C:227:HOH:O	2.08	0.53
1:A:130:MET:CG	1:B:138:GLN:HG3	2.38	0.53
1:B:96:HIS:HB2	4:B:243:HOH:O	2.08	0.53
1:B:175:SER:HB3	1:B:178:GLU:HG3	1.91	0.52
1:C:198:PHE:O	1:C:199:LYS:HG3	2.08	0.52
1:D:43:TRP:CE2	1:D:44:LEU:HG	2.44	0.52
2:B:225:C2E:N11	3:B:226:TAR:O11	2.28	0.52
1:A:198:PHE:O	1:A:199:LYS:HG3	2.10	0.52
1:D:37:THR:HG21	1:D:45:GLU:HG3	1.92	0.52
1:C:200:LYS:O	1:C:201:ILE:C	2.49	0.52
1:B:165:GLN:O	1:B:169:LEU:HG	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:TYR:C	1:A:156:TYR:CD2	2.84	0.51
1:D:212:TRP:O	1:D:216:ASN:HB2	2.10	0.51
1:A:200:LYS:O	1:A:201:ILE:C	2.46	0.51
1:C:130:MET:HG3	1:D:138:GLN:HG3	1.91	0.51
1:C:185:VAL:CG1	1:C:189:THR:HB	2.41	0.51
1:B:164:GLN:O	1:B:167:ILE:HG22	2.11	0.51
1:C:143:HIS:HD2	1:C:144:TYR:CD1	2.28	0.51
1:A:5:ASN:C	1:A:5:ASN:ND2	2.64	0.51
1:B:185:VAL:HG12	1:B:189:THR:HB	1.93	0.51
1:B:209:ALA:O	1:B:212:TRP:HB3	2.11	0.50
1:C:18:GLN:H	1:C:18:GLN:NE2	2.05	0.50
1:B:127:GLN:O	1:B:128:ASP:HB2	2.11	0.50
1:D:132:LEU:HD22	1:D:136:LEU:HD23	1.93	0.50
1:D:194:LEU:O	1:D:198:PHE:CD1	2.65	0.50
1:D:207:LEU:O	1:D:211:ILE:HG13	2.13	0.49
1:A:185:VAL:CG1	1:A:189:THR:HB	2.42	0.49
1:C:133:THR:HG23	1:C:136:LEU:H	1.77	0.49
1:D:74:SER:O	1:D:78:ILE:HG12	2.13	0.49
1:A:156:TYR:CE1	1:A:217:ILE:HD13	2.47	0.49
1:C:18:GLN:HE21	1:C:18:GLN:N	2.04	0.49
1:A:61:TYR:O	1:A:64:ILE:HG22	2.13	0.48
1:A:171:GLY:HA3	1:A:213:ALA:HB1	1.95	0.48
1:B:185:VAL:HG12	1:B:186:SER:N	2.27	0.48
1:A:185:VAL:HG13	1:A:189:THR:HB	1.95	0.48
1:C:144:TYR:HE2	1:D:145:ARG:HD3	1.79	0.48
1:B:167:ILE:CG2	1:B:168:LYS:N	2.76	0.48
1:D:96:HIS:CD2	1:D:139:GLU:HG3	2.48	0.48
1:D:210:LEU:HD23	1:D:210:LEU:N	2.21	0.48
1:A:112:ASP:OD1	1:A:112:ASP:N	2.45	0.48
1:D:207:LEU:HA	1:D:210:LEU:CD2	2.43	0.48
1:A:16:CYS:O	1:A:17:MET:C	2.53	0.47
1:D:215:ASN:O	1:D:216:ASN:CB	2.63	0.47
1:B:4:GLU:O	1:B:5:ASN:HB3	2.14	0.47
1:A:200:LYS:C	1:A:202:ASN:N	2.66	0.47
1:B:50:GLU:H	1:B:50:GLU:CD	2.18	0.47
1:C:148:ASN:N	1:C:148:ASN:ND2	2.62	0.47
1:C:185:VAL:HG13	1:C:189:THR:CB	2.45	0.47
1:B:171:GLY:HA3	1:B:213:ALA:HB1	1.96	0.47
1:D:109:PHE:CE1	1:D:131:TRP:CD1	3.03	0.46
1:D:127:GLN:O	1:D:128:ASP:HB2	2.14	0.46
1:D:143:HIS:HD2	1:D:144:TYR:CD2	2.34	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:ALA:HB1	1:A:205:ASN:O	2.16	0.46
1:D:215:ASN:O	1:D:216:ASN:CG	2.53	0.46
1:A:3:ASP:HB3	1:A:6:LYS:HD2	1.98	0.46
1:B:95:GLU:HB3	1:B:98:LEU:HG	1.97	0.46
1:A:72:TYR:CE2	1:A:102:TRP:HZ3	2.33	0.46
1:A:187:GLU:O	1:A:190:VAL:HG13	2.16	0.46
1:C:130:MET:SD	1:D:138:GLN:HG3	2.56	0.46
1:C:201:ILE:O	1:C:202:ASN:CB	2.64	0.45
1:D:41:GLU:O	1:D:41:GLU:HG2	2.15	0.45
1:B:180:ALA:HB1	1:B:185:VAL:O	2.16	0.45
1:D:206:ARG:O	1:D:209:ALA:HB3	2.17	0.45
1:A:197:VAL:C	1:A:199:LYS:N	2.69	0.45
1:D:50:GLU:H	1:D:50:GLU:CD	2.18	0.45
1:D:23:LYS:HE3	1:D:34:LEU:O	2.16	0.45
1:A:197:VAL:O	1:A:201:ILE:HG23	2.17	0.45
1:B:156:TYR:CD1	1:B:217:ILE:HG12	2.52	0.45
1:B:156:TYR:CZ	1:B:217:ILE:HD11	2.52	0.45
1:B:4:GLU:C	1:B:6:LYS:H	2.21	0.44
1:A:14:ASP:OD2	1:A:63:ARG:HD2	2.17	0.44
1:B:10:ARG:HB3	1:B:57:LEU:HD23	1.99	0.44
1:A:42:LEU:HA	1:A:42:LEU:HD12	1.72	0.44
1:C:187:GLU:O	1:C:190:VAL:HG13	2.17	0.44
1:A:3:ASP:OD2	1:A:5:ASN:HB3	2.18	0.44
1:A:122:MET:HB3	1:A:122:MET:HE2	1.73	0.44
1:B:177:ILE:HD11	1:B:187:GLU:OE1	2.17	0.44
1:A:105:LEU:HD23	1:A:140:TYR:OH	2.18	0.44
1:A:197:VAL:O	1:A:199:LYS:N	2.50	0.44
1:A:198:PHE:C	1:A:199:LYS:HG3	2.38	0.44
1:C:156:TYR:C	1:C:156:TYR:CD2	2.90	0.44
1:C:197:VAL:C	1:C:199:LYS:H	2.21	0.44
1:C:200:LYS:O	1:C:200:LYS:HG2	2.18	0.43
1:A:201:ILE:O	1:A:202:ASN:CB	2.66	0.43
1:A:130:MET:SD	1:B:138:GLN:HG3	2.58	0.43
1:B:74:SER:O	1:B:78:ILE:HG12	2.17	0.43
1:C:87:ILE:HD11	1:C:108:VAL:HG13	1.99	0.43
1:C:167:ILE:HD12	1:C:167:ILE:HA	1.77	0.43
1:D:9:VAL:HG21	1:D:32:LEU:HD22	2.00	0.43
1:A:130:MET:CE	1:A:132:LEU:HD11	2.48	0.43
1:B:159:LEU:HD12	1:B:164:GLN:HG3	2.00	0.43
1:A:164:GLN:O	1:A:167:ILE:HG22	2.19	0.43
1:A:167:ILE:HG23	1:A:168:LYS:N	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:4:GLU:CG	1:B:5:ASN:H	2.23	0.43
1:B:167:ILE:HD12	1:B:167:ILE:HA	1.85	0.43
1:B:156:TYR:OH	1:B:168:LYS:HD3	2.18	0.43
1:B:185:VAL:CG1	1:B:189:THR:HB	2.49	0.43
1:B:43:TRP:CZ2	1:B:44:LEU:CD2	3.02	0.43
1:A:87:ILE:HD13	1:A:105:LEU:HD11	1.99	0.42
1:A:167:ILE:O	1:A:168:LYS:C	2.57	0.42
1:B:167:ILE:HG23	1:B:168:LYS:N	2.34	0.42
1:C:167:ILE:CG2	1:C:168:LYS:N	2.82	0.42
1:B:159:LEU:HD12	1:B:164:GLN:CG	2.49	0.42
1:A:201:ILE:HD11	1:A:203:ALA:HB2	2.02	0.42
1:B:145:ARG:NH1	1:B:145:ARG:CG	2.69	0.42
1:D:133:THR:HG22	1:D:135:LYS:H	1.84	0.42
1:A:45:GLU:HB2	1:A:75:PHE:CE1	2.55	0.42
1:A:133:THR:HG22	1:A:136:LEU:N	2.26	0.42
1:D:49:PRO:HA	1:D:52:ARG:NH1	2.35	0.42
1:A:65:SER:O	1:A:68:VAL:HG22	2.20	0.41
1:D:133:THR:CG2	1:D:135:LYS:H	2.33	0.41
1:A:156:TYR:OH	1:A:168:LYS:CE	2.64	0.41
1:C:162:ARG:HA	1:C:162:ARG:HD3	1.90	0.41
1:A:185:VAL:CG1	1:A:186:SER:N	2.84	0.41
1:B:12:LEU:HA	1:B:37:THR:O	2.20	0.41
1:B:183:LEU:O	1:B:185:VAL:HG23	2.20	0.41
1:C:164:GLN:O	1:C:167:ILE:HG22	2.20	0.41
1:C:165:GLN:O	1:C:169:LEU:HG	2.21	0.41
1:A:154:GLN:NE2	1:A:154:GLN:HA	2.36	0.41
1:A:170:LEU:HD23	1:A:170:LEU:HA	1.92	0.41
1:D:143:HIS:HD2	1:D:144:TYR:CE2	2.39	0.41
1:A:45:GLU:HB2	1:A:75:PHE:HE1	1.86	0.41
1:A:167:ILE:HD12	1:A:167:ILE:HA	1.98	0.41
1:A:185:VAL:HG13	1:A:189:THR:CB	2.51	0.41
1:A:218:GLY:O	1:A:219:ILE:C	2.60	0.41
1:B:16:CYS:O	1:B:17:MET:C	2.59	0.41
2:B:225:C2E:HN11	3:B:226:TAR:HO1	1.59	0.41
1:D:4:GLU:CG	1:D:5:ASN:N	2.82	0.41
1:D:108:VAL:HG23	1:D:132:LEU:HD23	2.02	0.41
1:D:131:TRP:CH2	2:D:225:C2E:C41	3.04	0.41
1:A:72:TYR:OH	1:A:85:GLU:OE1	2.35	0.41
1:B:43:TRP:CZ2	1:B:44:LEU:HD22	2.56	0.41
1:B:185:VAL:CG1	1:B:186:SER:N	2.84	0.40
1:C:145:ARG:CG	1:C:145:ARG:NH1	2.68	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:206:ARG:O	1:D:210:LEU:HD23	2.20	0.40
1:B:122:MET:HB3	1:B:122:MET:HE2	1.82	0.40
1:B:161:LYS:HD2	1:B:161:LYS:H	1.87	0.40
1:B:161:LYS:HD2	1:B:161:LYS:N	2.36	0.40
1:B:217:ILE:HD12	1:B:217:ILE:O	2.22	0.40
1:B:177:ILE:HD12	1:B:177:ILE:N	2.36	0.40
1:A:200:LYS:O	1:A:202:ASN:N	2.54	0.40
1:B:39:PHE:O	1:B:42:LEU:HB2	2.22	0.40
1:D:126:LEU:HA	1:D:126:LEU:HD12	1.74	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	215/225 (96%)	201 (94%)	12 (6%)	2 (1%)	17 46
1	B	214/225 (95%)	202 (94%)	11 (5%)	1 (0%)	29 61
1	C	211/225 (94%)	197 (93%)	14 (7%)	0	100 100
1	D	169/225 (75%)	156 (92%)	13 (8%)	0	100 100
All	All	809/900 (90%)	756 (93%)	50 (6%)	3 (0%)	34 66

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	198	PHE
1	B	202	ASN
1	A	143	HIS

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	199/207 (96%)	175 (88%)	24 (12%)	5	15
1	B	198/207 (96%)	176 (89%)	22 (11%)	6	19
1	C	196/207 (95%)	170 (87%)	26 (13%)	4	12
1	D	161/207 (78%)	140 (87%)	21 (13%)	4	13
All	All	754/828 (91%)	661 (88%)	93 (12%)	4	15

All (93) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	10	ARG
1	A	12	LEU
1	A	15	VAL
1	A	18	GLN
1	A	29	LYS
1	A	50	GLU
1	A	61	TYR
1	A	85	GLU
1	A	94	ILE
1	A	116	ASP
1	A	117	THR
1	A	133	THR
1	A	142	LEU
1	A	151	VAL
1	A	160	THR
1	A	161	LYS
1	A	170	LEU
1	A	185	VAL
1	A	190	VAL
1	A	199	LYS
1	A	201	ILE
1	A	202	ASN
1	A	217	ILE

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Mol	Chain	Res	Type
1	B	12	LEU
1	B	18	GLN
1	B	29	LYS
1	B	35	GLU
1	B	44	LEU
1	B	61	TYR
1	B	70	THR
1	B	71	ASP
1	B	76	LYS
1	B	101	LYS
1	B	111	ILE
1	B	112	ASP
1	B	133	THR
1	B	142	LEU
1	B	145	ARG
1	B	151	VAL
1	B	159	LEU
1	B	170	LEU
1	B	190	VAL
1	B	202	ASN
1	B	210	LEU
1	B	217	ILE
1	C	10	ARG
1	C	15	VAL
1	C	17	MET
1	C	18	GLN
1	C	29	LYS
1	C	35	GLU
1	C	61	TYR
1	C	76	LYS
1	C	87	ILE
1	C	93	ASP
1	C	94	ILE
1	C	117	THR
1	C	139	GLU
1	C	142	LEU
1	C	145	ARG
1	C	148	ASN
1	C	159	LEU
1	C	161	LYS
1	C	170	LEU
1	C	185	VAL

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Mol	Chain	Res	Type
1	C	187	GLU
1	C	190	VAL
1	C	199	LYS
1	C	201	ILE
1	C	202	ASN
1	C	210	LEU
1	D	10	ARG
1	D	15	VAL
1	D	18	GLN
1	D	29	LYS
1	D	32	LEU
1	D	34	LEU
1	D	35	GLU
1	D	50	GLU
1	D	55	GLN
1	D	71	ASP
1	D	82	ASP
1	D	85	GLU
1	D	92	GLN
1	D	126	LEU
1	D	133	THR
1	D	145	ARG
1	D	152	THR
1	D	193	HIS
1	D	197	VAL
1	D	199	LYS
1	D	210	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	18	GLN
1	A	92	GLN
1	A	127	GLN
1	A	154	GLN
1	B	18	GLN
1	B	103	ASN
1	B	143	HIS
1	C	18	GLN
1	C	89	ASN
1	C	96	HIS

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Mol	Chain	Res	Type
1	C	103	ASN
1	C	143	HIS
1	C	148	ASN
1	C	154	GLN
1	C	193	HIS
1	C	196	ASN
1	D	18	GLN
1	D	89	ASN
1	D	92	GLN
1	D	143	HIS
1	D	148	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

6 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	C2E	C	225	-	44,52,52	1.11	6 (13%)	52,82,82	1.18	6 (11%)
2	C2E	A	225	-	44,52,52	1.18	6 (13%)	52,82,82	1.35	9 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	C2E	D	225	-	44,52,52	1.14	5 (11%)	52,82,82	1.24	6 (11%)
2	C2E	B	225	-	44,52,52	1.23	5 (11%)	52,82,82	1.29	6 (11%)
3	TAR	B	226	-	9,9,9	1.86	2 (22%)	12,12,12	1.60	3 (25%)
3	TAR	D	226	-	9,9,9	1.86	2 (22%)	12,12,12	1.39	2 (16%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	C2E	C	225	-	-	5/22/62/62	0/6/7/7
2	C2E	A	225	-	-	0/22/62/62	0/6/7/7
2	C2E	D	225	-	-	3/22/62/62	0/6/7/7
2	C2E	B	225	-	-	2/22/62/62	0/6/7/7
3	TAR	B	226	-	-	5/12/12/12	-
3	TAR	D	226	-	2/2/4/4	3/12/12/12	-

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	226	TAR	O41-C4	3.92	1.43	1.30
3	D	226	TAR	O41-C4	3.62	1.42	1.30
3	D	226	TAR	O11-C1	3.62	1.42	1.30
3	B	226	TAR	O11-C1	3.46	1.41	1.30
2	A	225	C2E	C51-C61	-3.21	1.40	1.47
2	B	225	C2E	C5-C6	-2.94	1.41	1.47
2	B	225	C2E	C51-C61	-2.84	1.41	1.47
2	D	225	C2E	C81-N71	-2.82	1.30	1.35
2	A	225	C2E	C5-C6	-2.75	1.41	1.47
2	D	225	C2E	C5-C6	-2.71	1.41	1.47
2	C	225	C2E	C51-C61	-2.68	1.42	1.47
2	B	225	C2E	C81-N71	-2.66	1.30	1.35
2	D	225	C2E	C8-N7	-2.63	1.30	1.35
2	B	225	C2E	C8-N7	-2.52	1.30	1.35
2	B	225	C2E	C61-N11	-2.51	1.34	1.37
2	A	225	C2E	C81-N71	-2.45	1.30	1.35
2	C	225	C2E	O4A-C1A	2.36	1.44	1.41
2	C	225	C2E	C81-N71	-2.31	1.31	1.35
2	C	225	C2E	O4'-C1'	2.22	1.44	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	225	C2E	C8-N7	-2.22	1.31	1.35
2	D	225	C2E	O4A-C1A	2.21	1.44	1.41
2	C	225	C2E	C8-N7	-2.19	1.31	1.35
2	C	225	C2E	C5-C6	-2.14	1.43	1.47
2	D	225	C2E	O4'-C1'	2.11	1.44	1.41
2	A	225	C2E	C6-N1	-2.05	1.34	1.37
2	A	225	C2E	C61-N11	-2.05	1.34	1.37

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	225	C2E	C81-N71-C51	3.71	110.05	102.99
2	D	225	C2E	C21-N11-C61	-3.30	119.02	125.10
2	C	225	C2E	C81-N71-C51	3.30	109.28	102.99
2	A	225	C2E	C8-N7-C5	3.27	109.22	102.99
2	B	225	C2E	C8-N7-C5	3.26	109.21	102.99
2	C	225	C2E	C8-N7-C5	3.24	109.16	102.99
2	B	225	C2E	C2-N1-C6	-3.13	119.34	125.10
2	B	225	C2E	C5-C6-N1	3.09	119.41	113.95
2	D	225	C2E	C8-N7-C5	3.08	108.86	102.99
2	A	225	C2E	C21-N11-C61	-3.03	119.53	125.10
2	A	225	C2E	C81-N71-C51	2.98	108.66	102.99
3	B	226	TAR	O11-C1-O1	-2.87	117.56	124.09
2	A	225	C2E	C5-C6-N1	2.85	118.98	113.95
2	B	225	C2E	C81-N71-C51	2.84	108.40	102.99
2	A	225	C2E	C51-C61-N11	2.81	118.91	113.95
2	C	225	C2E	C21-N11-C61	-2.79	119.96	125.10
2	B	225	C2E	C51-C61-N11	2.76	118.83	113.95
2	D	225	C2E	C2-N1-C6	-2.75	120.03	125.10
2	D	225	C2E	C5-C6-N1	2.73	118.77	113.95
2	A	225	C2E	C2-N1-C6	-2.71	120.11	125.10
2	C	225	C2E	C51-C61-N11	2.65	118.62	113.95
2	D	225	C2E	C51-C61-N11	2.64	118.62	113.95
3	B	226	TAR	O11-C1-C2	2.51	120.05	113.27
3	D	226	TAR	O3-C3-C4	2.50	115.90	110.66
2	A	225	C2E	C2A-C3A-C4A	-2.45	98.89	103.22
2	B	225	C2E	C21-N11-C61	-2.42	120.64	125.10
2	C	225	C2E	C5-C6-N1	2.25	117.92	113.95
2	A	225	C2E	C5'-C4'-C3'	-2.16	107.24	114.40
2	C	225	C2E	C2-N1-C6	-2.14	121.16	125.10
3	D	226	TAR	O3-C3-C2	-2.06	106.14	110.23
3	B	226	TAR	O1-C1-C2	-2.03	116.29	121.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	225	C2E	O61-C61-C51	-2.02	120.43	124.37

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	D	226	TAR	C2
3	D	226	TAR	C3

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	226	TAR	O1-C1-C2-O2
3	B	226	TAR	O11-C1-C2-O2
3	B	226	TAR	O3-C3-C4-O4
3	B	226	TAR	O3-C3-C4-O41
2	C	225	C2E	O4A-C4A-C5A-O5A
2	C	225	C2E	C3A-C4A-C5A-O5A
3	B	226	TAR	O11-C1-C2-C3
3	D	226	TAR	O11-C1-C2-C3
2	C	225	C2E	C3A-O3A-P1-O5'
2	B	225	C2E	C4A-C3A-O3A-P1
3	D	226	TAR	O11-C1-C2-O2
2	D	225	C2E	C4A-C3A-O3A-P1
2	B	225	C2E	O4'-C4'-C5'-O5'
2	D	225	C2E	C2A-C3A-O3A-P1
2	C	225	C2E	C5A-O5A-P11-O3'
2	C	225	C2E	O4'-C4'-C5'-O5'
3	D	226	TAR	O1-C1-C2-C3
2	D	225	C2E	C2'-C3'-O3'-P11

There are no ring outliers.

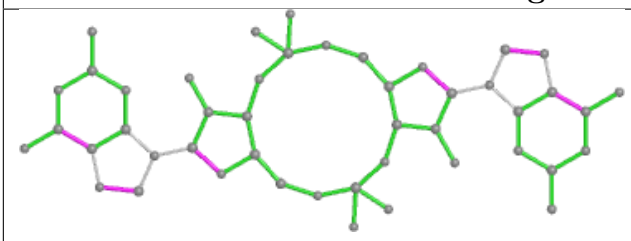
4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	225	C2E	1	0
2	D	225	C2E	1	0
2	B	225	C2E	2	0
3	B	226	TAR	4	0

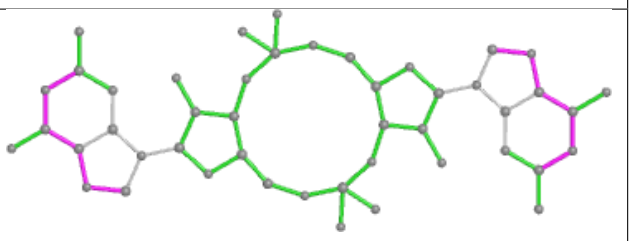
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

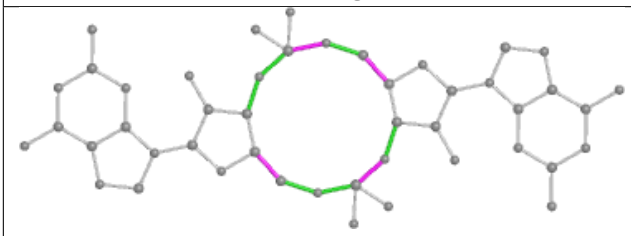
Ligand C2E C 225



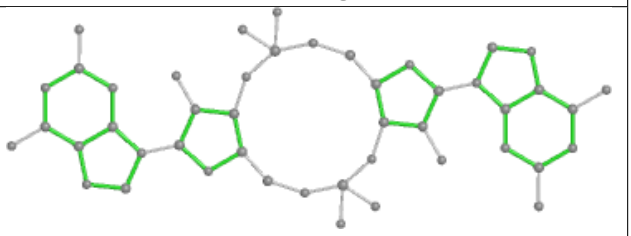
Bond lengths



Bond angles

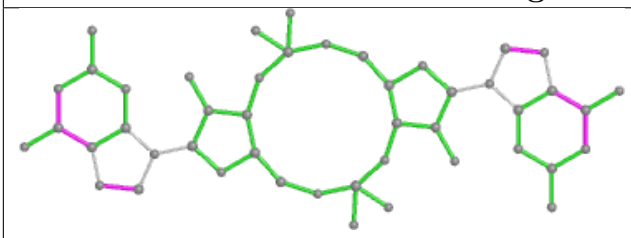


Torsions

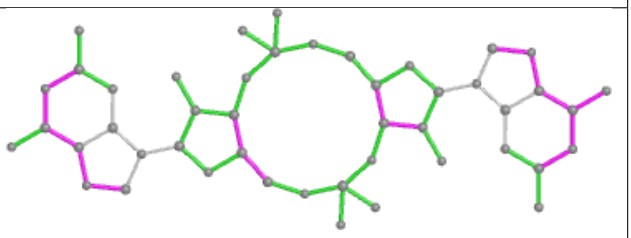


Rings

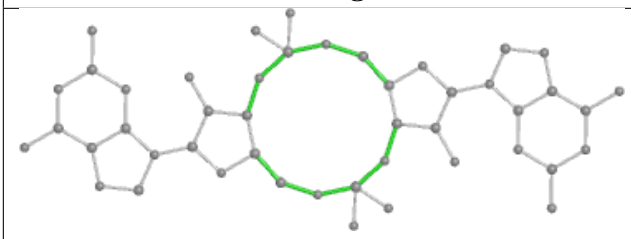
Ligand C2E A 225



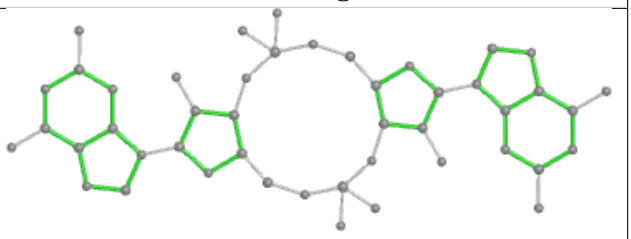
Bond lengths



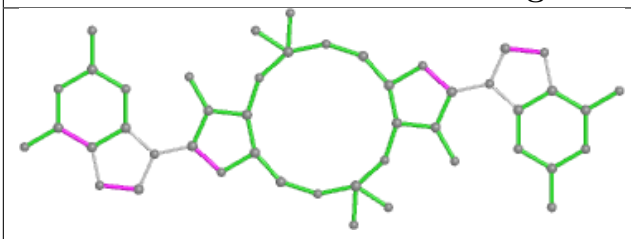
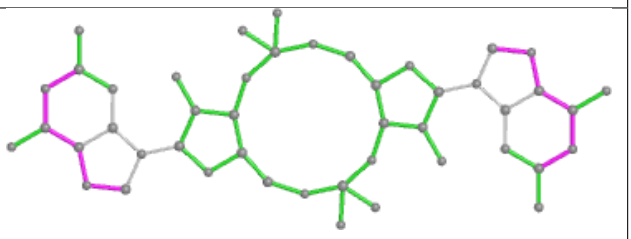
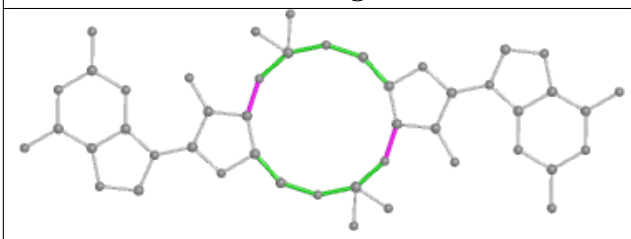
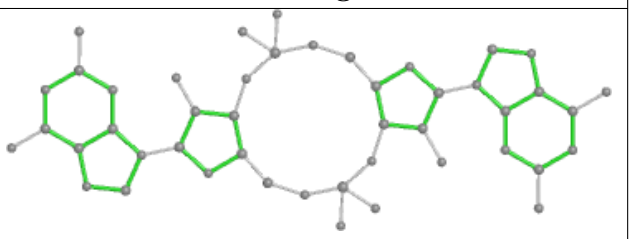
Bond angles

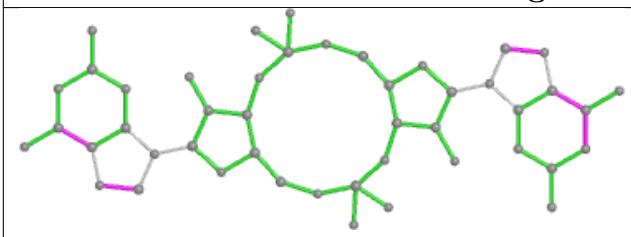
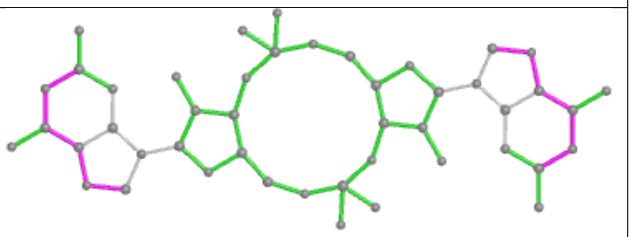
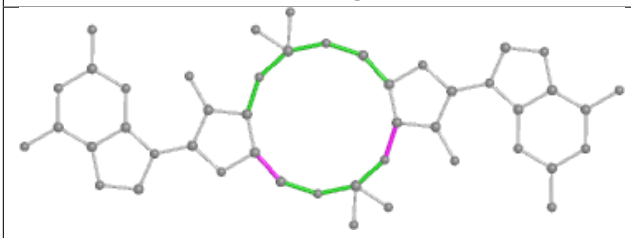
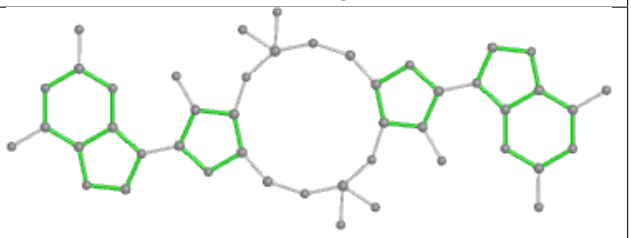


Torsions



Rings

Ligand C2E D 225	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand C2E B 225	
 Bond lengths	 Bond angles
 Torsions	 Rings

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	217/225 (96%)	0.02	1 (0%)	91	88	18, 41, 72, 96	0
1	B	216/225 (96%)	0.23	9 (4%)	36	26	20, 41, 82, 97	0
1	C	213/225 (94%)	0.60	27 (12%)	3	2	31, 62, 98, 117	0
1	D	173/225 (76%)	0.93	32 (18%)	1	1	41, 61, 107, 125	0
All	All	819/900 (91%)	0.42	69 (8%)	11	5	18, 52, 95, 125	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	149	SER	8.2
1	D	210	LEU	7.2
1	D	215	ASN	6.7
1	D	197	VAL	6.7
1	D	201	ILE	6.1
1	D	195	HIS	6.0
1	D	211	ILE	6.0
1	D	151	VAL	5.9
1	D	202	ASN	5.7
1	D	203	ALA	5.1
1	C	215	ASN	5.0
1	D	198	PHE	5.0
1	D	207	LEU	5.0
1	D	147	GLY	4.9
1	C	155	MET	4.9
1	A	3	ASP	4.7
1	C	214	LYS	4.5
1	B	188	ASN	4.1
1	D	121	GLY	3.8
1	B	172	SER	3.8
1	D	150	VAL	3.7

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Mol	Chain	Res	Type	RSRZ
1	D	194	LEU	3.7
1	D	49	PRO	3.6
1	B	217	ILE	3.4
1	C	159	LEU	3.4
1	D	8	ASN	3.4
1	D	148	ASN	3.3
1	B	171	GLY	3.3
1	D	208	GLN	3.2
1	C	158	LYS	3.2
1	C	204	LYS	3.2
1	C	154	GLN	3.2
1	C	144	TYR	3.2
1	C	200	LYS	3.1
1	C	150	VAL	3.1
1	C	211	ILE	3.1
1	D	205	ASN	3.1
1	C	47	ASN	3.0
1	D	214	LYS	2.9
1	D	55	GLN	2.9
1	B	181	ASP	2.8
1	D	200	LYS	2.7
1	B	179	ILE	2.7
1	C	210	LEU	2.6
1	D	149	SER	2.6
1	B	190	VAL	2.5
1	D	206	ARG	2.5
1	D	212	TRP	2.5
1	C	79	SER	2.5
1	D	204	LYS	2.4
1	C	52	ARG	2.3
1	D	44	LEU	2.3
1	C	50	GLU	2.3
1	C	197	VAL	2.3
1	C	4	GLU	2.3
1	C	156	TYR	2.3
1	C	151	VAL	2.3
1	D	152	THR	2.2
1	C	184	PHE	2.2
1	C	147	GLY	2.2
1	C	171	GLY	2.2
1	C	146	ALA	2.2
1	D	47	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	165	GLN	2.2
1	B	4	GLU	2.2
1	C	196	ASN	2.1
1	D	213	ALA	2.1
1	D	196	ASN	2.1
1	B	156	TYR	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 monosaccharides in this entry.

6.4 Ligands (i)

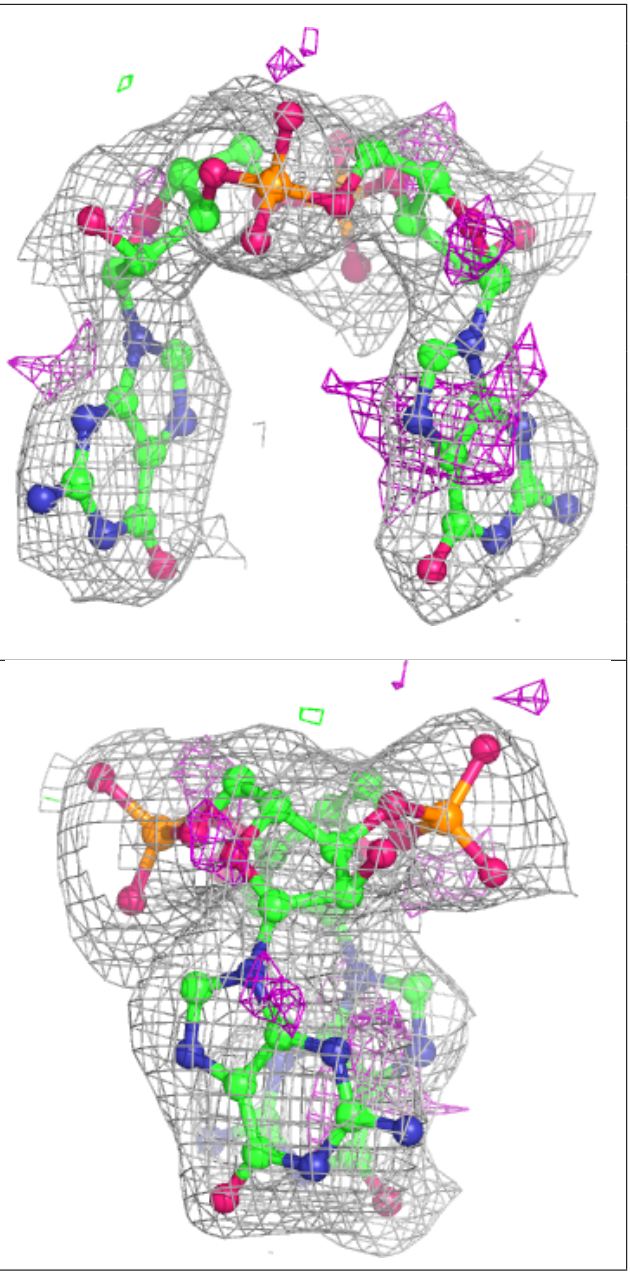
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	TAR	D	226	10/10	0.66	0.22	53,61,69,72	0
3	TAR	B	226	10/10	0.72	0.54	46,54,65,65	0
2	C2E	C	225	46/46	0.95	0.18	38,49,56,59	0
2	C2E	D	225	46/46	0.95	0.17	40,49,53,54	0
2	C2E	A	225	46/46	0.98	0.17	22,28,33,35	0
2	C2E	B	225	46/46	0.98	0.18	22,26,33,35	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.

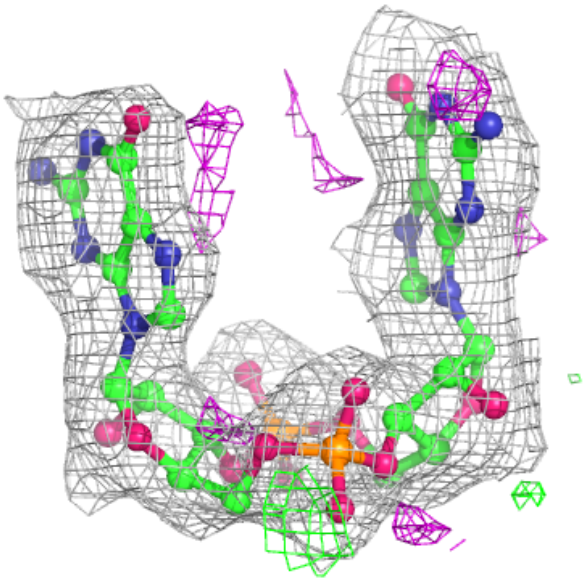
Electron density around C2E C 225:

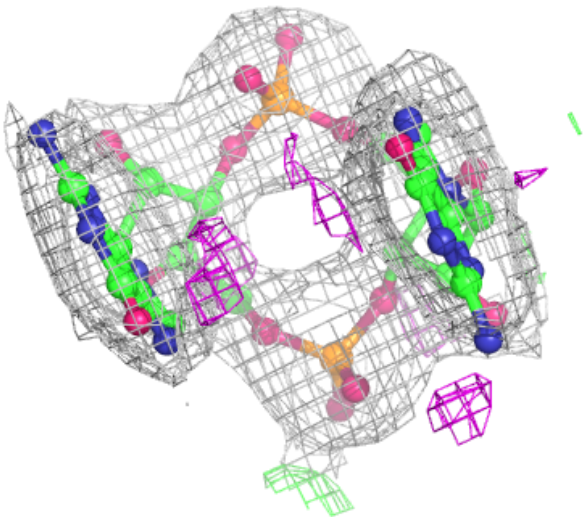
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

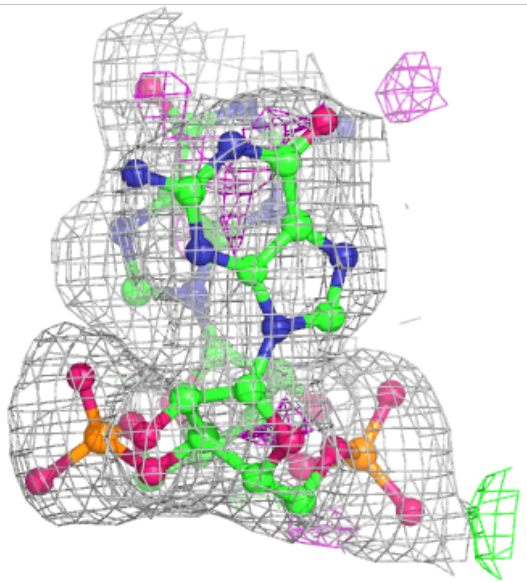


Electron density around C2E D 225:

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

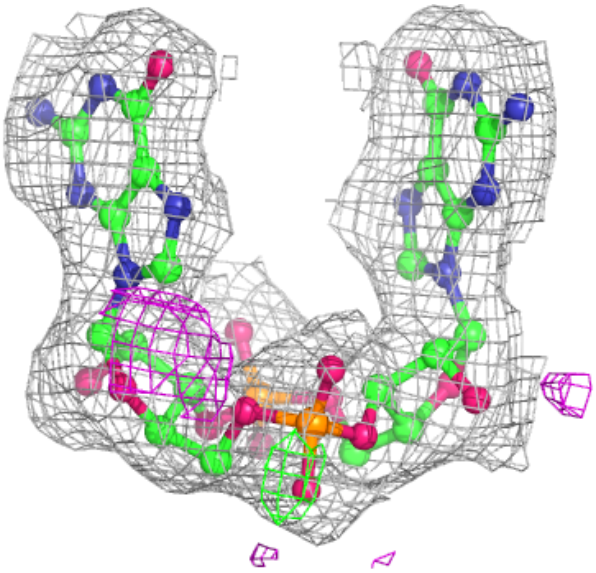


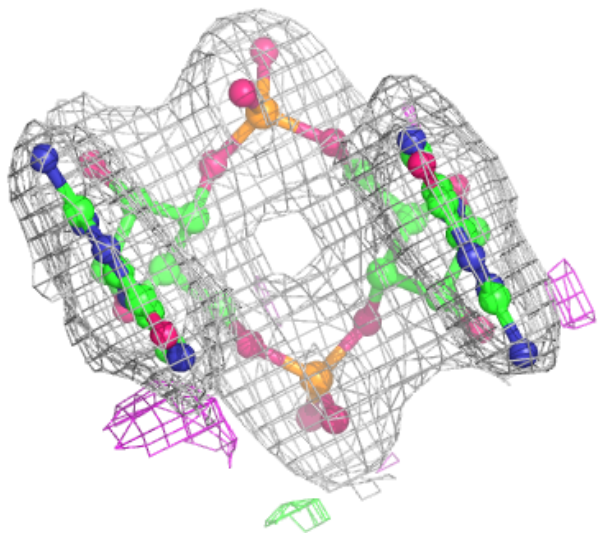


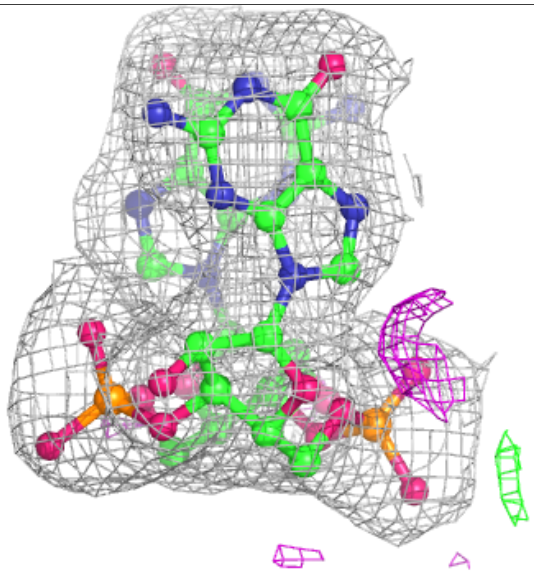


Electron density around C2E A 225:

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

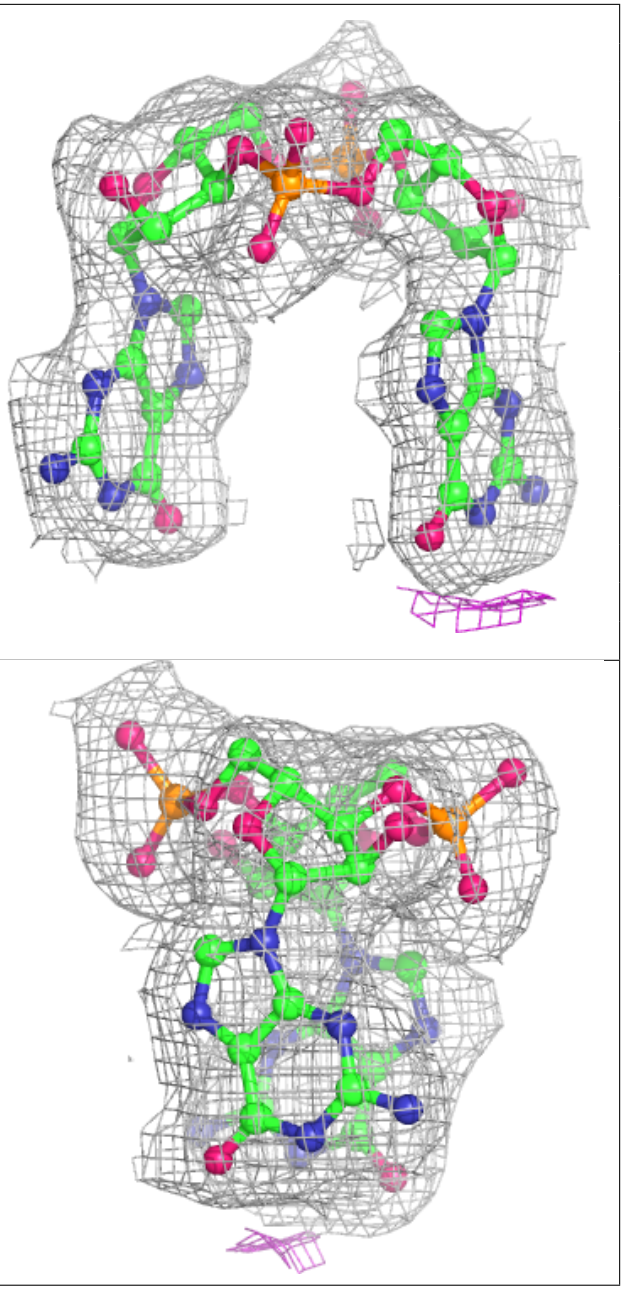






Electron density around C2E B 225:

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.