Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 06:24 pm BST

PDB ID	:	3E5X
Title	:	OCPA complexed CprK
Authors	:	Levy, C.
Deposited on	:	2008-08-14
Resolution	:	2.00 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:

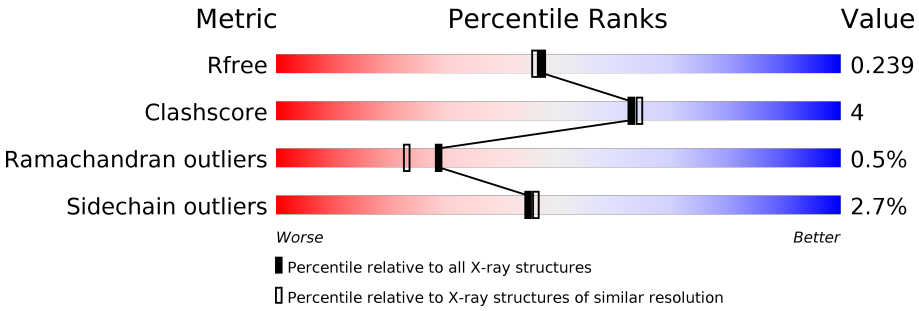
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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 <=5%

Mol	Chain	Length	Quality of chain
1	A	250	74% 10% • 14%
1	B	250	74% 12% • 12%
1	C	250	75% 9% • 14%
1	D	250	72% 14% • 12%

2 Entry composition

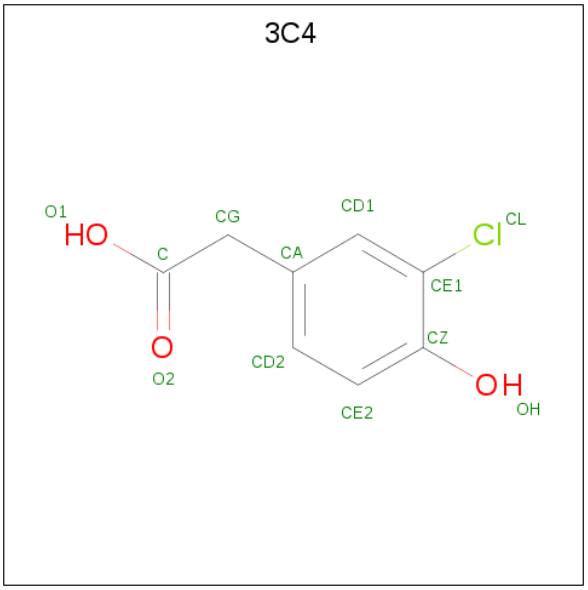
There are 3 unique types of molecules in this entry. The entry contains 7584 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Cyclic nucleotide-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	0	0
			1719	1108	283	317	11			
1	C	214	Total	C	N	O	S	0	0	0
			1719	1108	283	317	11			
1	B	219	Total	C	N	O	S	0	0	0
			1758	1133	291	323	11			
1	D	220	Total	C	N	O	S	0	1	0
			1775	1143	294	327	11			

- Molecule 2 is (3-CHLORO-4-HYDROXYPHENYL)ACETIC ACID (three-letter code: 3C4) (formula: $C_8H_7ClO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Cl	O	0	0
			12	8	1	3		
2	C	1	Total	C	Cl	O	0	0
			12	8	1	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	Cl	O	0	0
			12	8	1	3		
2	B	1	Total	C	Cl	O	0	0
			12	8	1	3		
2	D	1	Total	C	Cl	O	0	0
			12	8	1	3		

- Molecule 3 is water.

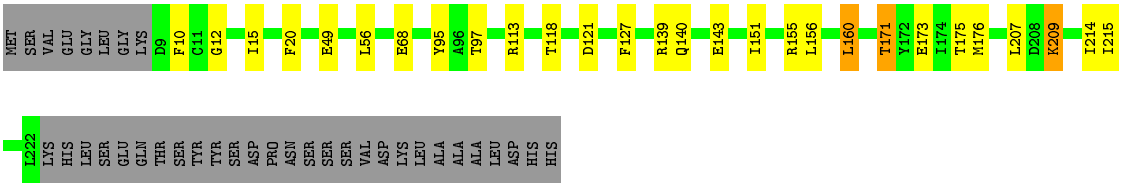
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	121	Total	O	0	0
			121	121		
3	C	130	Total	O	0	0
			130	130		
3	B	151	Total	O	0	0
			151	151		
3	D	151	Total	O	0	0
			151	151		

3 Residue-property plots (i)

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. 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. 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.

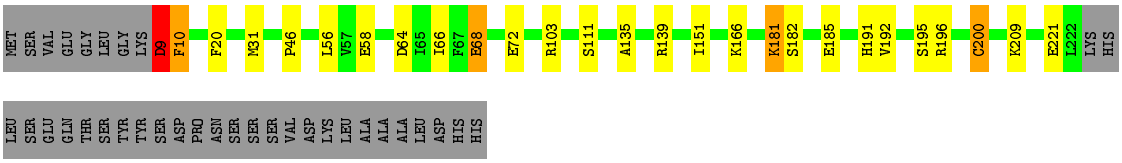
- Molecule 1: Cyclic nucleotide-binding protein

Chain A:



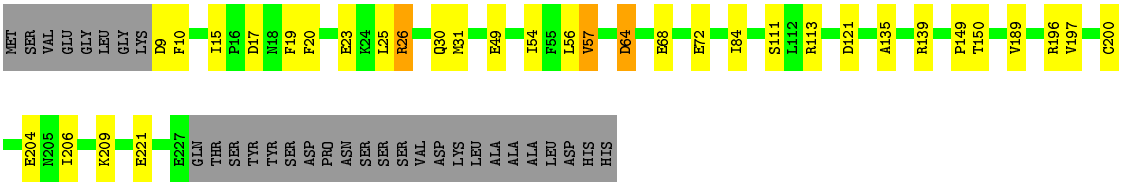
- Molecule 1: Cyclic nucleotide-binding protein

Chain C:



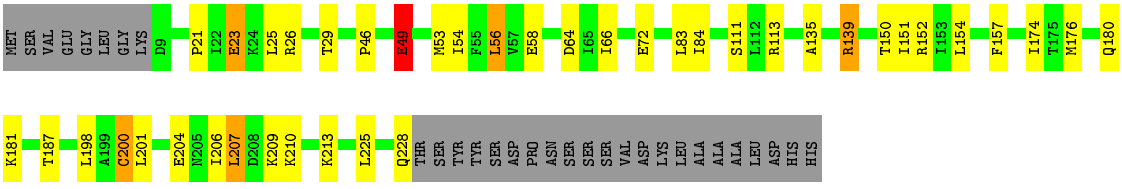
- Molecule 1: Cyclic nucleotide-binding protein

Chain B:



- Molecule 1: Cyclic nucleotide-binding protein

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	56.70Å 118.38Å 87.46Å 90.00° 95.68° 90.00°	Depositor
Resolution (Å)	35.94 – 2.00 32.92 – 2.00	Depositor EDS
% Data completeness (in resolution range)	96.3 (35.94-2.00) 96.3 (32.92-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.41 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.193 , 0.238 0.196 , 0.239	Depositor DCC
R_{free} test set	3754 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	28.9	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 36.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	7584	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% of the height of the origin peak. No significant pseudotranslation is detected.*

[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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 3C4

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	1.12	4/1752 (0.2%)	0.91	7/2362 (0.3%)
1	B	1.16	7/1792 (0.4%)	0.97	5/2415 (0.2%)
1	C	1.13	5/1752 (0.3%)	0.99	3/2362 (0.1%)
1	D	1.14	7/1813 (0.4%)	0.93	5/2444 (0.2%)
All	All	1.14	23/7109 (0.3%)	0.95	20/9583 (0.2%)

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	B	0	1
1	C	0	1
All	All	0	2

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	171	THR	CB-OG1	17.68	1.78	1.43
1	C	200	CYS	CB-SG	-11.46	1.62	1.82
1	B	200	CYS	CB-SG	-10.38	1.64	1.82
1	B	221	GLU	C-N	10.38	1.57	1.34
1	A	171	THR	CB-CG2	10.35	1.86	1.52
1	C	68	GLU	CG-CD	8.32	1.64	1.51
1	D	200	CYS	CB-SG	-8.16	1.68	1.82
1	D	23	GLU	CG-CD	7.99	1.64	1.51
1	D	49	GLU	CG-CD	7.59	1.63	1.51
1	A	68	GLU	CG-CD	7.35	1.62	1.51
1	B	135	ALA	CA-CB	6.89	1.67	1.52
1	A	49	GLU	CG-CD	6.28	1.61	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	49	GLU	CD-OE1	6.15	1.32	1.25
1	D	58	GLU	CG-CD	5.86	1.60	1.51
1	B	68	GLU	CG-CD	5.75	1.60	1.51
1	C	68	GLU	CB-CG	5.62	1.62	1.52
1	B	23	GLU	CG-CD	5.58	1.60	1.51
1	D	111	SER	CB-OG	-5.47	1.35	1.42
1	B	49	GLU	CG-CD	5.21	1.59	1.51
1	B	57	VAL	CB-CG2	5.13	1.63	1.52
1	D	135	ALA	CA-CB	5.12	1.63	1.52
1	C	135	ALA	CA-CB	5.09	1.63	1.52
1	C	195	SER	CB-OG	5.01	1.48	1.42

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	139	ARG	NE-CZ-NH2	-8.93	115.84	120.30
1	C	139	ARG	NE-CZ-NH1	8.81	124.70	120.30
1	A	139	ARG	NE-CZ-NH1	8.12	124.36	120.30
1	D	113	ARG	NE-CZ-NH1	7.73	124.17	120.30
1	A	56	LEU	CB-CG-CD1	-7.18	98.80	111.00
1	B	26	ARG	NE-CZ-NH2	-7.11	116.74	120.30
1	D	207	LEU	CA-CB-CG	6.62	130.52	115.30
1	A	139	ARG	NE-CZ-NH2	-6.56	117.02	120.30
1	C	196	ARG	NE-CZ-NH1	6.52	123.56	120.30
1	D	56	LEU	CB-CG-CD2	-6.10	100.64	111.00
1	A	127	PHE	CB-CG-CD1	6.08	125.06	120.80
1	B	10	PHE	N-CA-C	-5.62	95.81	111.00
1	D	139	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	B	196	ARG	NE-CZ-NH2	-5.40	117.60	120.30
1	B	113	ARG	NE-CZ-NH1	5.31	122.95	120.30
1	B	64	ASP	CB-CG-OD2	-5.20	113.62	118.30
1	A	121	ASP	CB-CG-OD1	5.18	122.96	118.30
1	A	113	ARG	NE-CZ-NH1	5.07	122.84	120.30
1	D	139	ARG	NE-CZ-NH2	-5.06	117.77	120.30
1	A	155	ARG	NE-CZ-NH1	5.06	122.83	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	9	ASP	Peptide
1	C	9	ASP	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	1719	0	1748	13	0
1	B	1758	0	1788	12	0
1	C	1719	0	1748	14	0
1	D	1775	0	1798	23	0
2	A	12	0	6	0	0
2	B	24	0	12	0	0
2	C	12	0	6	0	0
2	D	12	0	6	0	0
3	A	121	0	0	2	0
3	B	151	0	0	3	0
3	C	130	0	0	4	0
3	D	151	0	0	6	0
All	All	7584	0	7112	61	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 4.

All (61) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:THR:CG2	1:A:171:THR:CB	1.86	1.50
1:A:171:THR:CB	1:A:171:THR:OG1	1.78	1.30
1:B:121:ASP:OD2	3:B:1594:HOH:O	1.90	0.89
1:D:23:GLU:O	1:D:26:ARG:HG2	1.78	0.82
1:C:151:ILE:HD12	3:C:1099:HOH:O	1.80	0.80
1:B:139:ARG:NH1	3:B:1318:HOH:O	2.21	0.72
1:D:152:ARG:NH1	1:D:187:THR:O	2.33	0.61
1:D:157:PHE:CZ	1:D:198:LEU:HD21	2.38	0.58
1:D:150:THR:HG21	3:D:1116:HOH:O	2.04	0.57
1:D:49:GLU:HG2	3:D:1305:HOH:O	2.06	0.56
1:B:56:LEU:HD12	3:B:1575:HOH:O	2.06	0.56
1:D:174:ILE:HG22	1:D:176:MET:HG2	1.89	0.54
1:C:9:ASP:N	1:C:9:ASP:OD1	2.36	0.54
1:A:173:GLU:HG2	1:A:215:ILE:HD12	1.89	0.53
1:D:204:GLU:HB2	1:D:206:ILE:HD12	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:209:LYS:O	1:D:213:LYS:O	2.26	0.53
1:A:151:ILE:HD12	3:A:1258:HOH:O	2.09	0.53
1:D:157:PHE:HZ	1:D:198:LEU:HD21	1.74	0.53
1:C:68:GLU:CD	1:C:68:GLU:H	2.13	0.52
1:D:154:LEU:CD2	1:D:201:LEU:HD21	2.40	0.52
1:C:58:GLU:HG2	1:C:103:ARG:HB3	1.92	0.51
1:B:19:PHE:O	1:B:20:PHE:C	2.49	0.51
1:C:182:SER:HA	1:C:185:GLU:HG2	1.93	0.51
1:D:46:PRO:HB3	1:D:66:ILE:HD12	1.93	0.51
1:A:171:THR:CG2	1:A:171:THR:CA	2.84	0.50
1:D:180:GLN:HG2	1:D:198:LEU:HD12	1.94	0.48
1:C:191:HIS:HD2	3:C:1133:HOH:O	1.96	0.47
1:D:228:GLN:HB2	3:D:1583:HOH:O	2.13	0.47
1:D:228:GLN:C	3:D:1130:HOH:O	2.53	0.47
1:D:200:CYS:O	1:D:204:GLU:HG3	2.15	0.46
1:D:206:ILE:HD13	1:D:225:LEU:CD1	2.46	0.45
1:A:209:LYS:HG2	1:A:214:ILE:CD1	2.46	0.45
1:A:15:ILE:HD11	1:A:156:LEU:HD11	1.98	0.45
1:B:26:ARG:NH2	1:B:57:VAL:O	2.50	0.44
1:B:31:MET:O	1:B:111:SER:HB2	2.16	0.44
1:B:150:THR:HG22	1:B:197:VAL:HG21	1.99	0.44
1:C:181:LYS:O	1:C:185:GLU:HG2	2.17	0.44
1:C:46:PRO:HB3	1:C:66:ILE:HD12	2.01	0.43
1:B:149:PRO:HB3	1:B:189:VAL:HG22	2.00	0.43
1:C:182:SER:OG	3:C:1534:HOH:O	2.21	0.43
1:D:139:ARG:CZ	3:D:1502:HOH:O	2.67	0.43
1:C:31:MET:O	1:C:111:SER:HB2	2.19	0.42
1:B:64:ASP:OD2	1:B:72:GLU:OE2	2.37	0.42
1:D:151:ILE:HD13	1:D:151:ILE:HA	1.86	0.42
1:C:64:ASP:OD2	1:C:72:GLU:OE2	2.37	0.42
1:B:15:ILE:HD12	1:B:17:ASP:HB3	2.00	0.42
1:D:53:MET:HB3	1:D:83:LEU:HD11	2.01	0.42
1:A:95:TYR:CZ	1:A:97:THR:CG2	3.02	0.42
1:B:204:GLU:HB2	1:B:206:ILE:HD12	2.01	0.42
1:D:150:THR:HG23	3:D:1208:HOH:O	2.19	0.41
1:C:192:VAL:HG12	3:C:1343:HOH:O	2.20	0.41
1:A:140:GLN:HG3	3:A:1263:HOH:O	2.21	0.41
1:D:26:ARG:HA	1:D:29:THR:HG23	2.03	0.41
1:B:54:ILE:HB	1:B:84:ILE:HB	2.03	0.41
1:A:10:PHE:CE2	1:A:12:GLY:HA3	2.56	0.41
1:A:118:THR:O	1:C:166:LYS:HB3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:221:GLU:HA	1:C:221:GLU:OE1	2.20	0.41
1:A:209:LYS:HE2	1:A:209:LYS:O	2.21	0.41
1:D:54:ILE:HB	1:D:84:ILE:HB	2.03	0.40
1:A:160:LEU:HA	1:A:160:LEU:HD12	1.95	0.40
1:D:64:ASP:OD2	1:D:72:GLU:OE2	2.38	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	212/250 (85%)	206 (97%)	5 (2%)	1 (0%)	29	23
1	B	217/250 (87%)	211 (97%)	6 (3%)	0	100	100
1	C	212/250 (85%)	204 (96%)	6 (3%)	2 (1%)	17	11
1	D	219/250 (88%)	214 (98%)	4 (2%)	1 (0%)	29	23
All	All	860/1000 (86%)	835 (97%)	21 (2%)	4 (0%)	29	23

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	210	LYS
1	C	10	PHE
1	C	20	PHE
1	A	20	PHE

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	189/220 (86%)	183 (97%)	6 (3%)	39	38
1	B	193/220 (88%)	190 (98%)	3 (2%)	62	67
1	C	189/220 (86%)	183 (97%)	6 (3%)	39	38
1	D	195/220 (89%)	189 (97%)	6 (3%)	40	40
All	All	766/880 (87%)	745 (97%)	21 (3%)	44	46

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	143	GLU
1	A	160	LEU
1	A	175	THR
1	A	176	MET
1	A	207	LEU
1	A	209	LYS
1	C	9	ASP
1	C	10	PHE
1	C	56	LEU
1	C	181	LYS
1	C	200	CYS
1	C	209	LYS
1	B	25	LEU
1	B	30	GLN
1	B	209	LYS
1	D	21	PRO
1	D	25	LEU
1	D	49	GLU
1	D	56	LEU
1	D	181	LYS
1	D	207	LEU

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

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

5 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	3C4	B	505	-	9,12,12	4.59	6 (66%)	12,16,16	3.67	6 (50%)
2	3C4	D	503	-	9,12,12	5.21	4 (44%)	12,16,16	2.52	3 (25%)
2	3C4	C	502	-	9,12,12	4.78	3 (33%)	12,16,16	2.37	3 (25%)
2	3C4	B	504	-	9,12,12	4.70	4 (44%)	12,16,16	2.52	3 (25%)
2	3C4	A	501	-	9,12,12	5.16	4 (44%)	12,16,16	2.70	6 (50%)

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	3C4	B	505	-	-	0/2/4/4	0/1/1/1
2	3C4	D	503	-	-	0/2/4/4	0/1/1/1
2	3C4	C	502	-	-	0/2/4/4	0/1/1/1
2	3C4	B	504	-	-	0/2/4/4	0/1/1/1
2	3C4	A	501	-	-	0/2/4/4	0/1/1/1

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	3C4	CZ-CE1	12.67	1.51	1.39
2	D	503	3C4	CZ-CE1	12.63	1.51	1.39
2	C	502	3C4	CZ-CE1	11.51	1.50	1.39
2	B	504	3C4	CZ-CE1	11.48	1.50	1.39
2	B	505	3C4	CZ-CE1	9.02	1.48	1.39
2	B	505	3C4	CE2-CD2	-7.33	1.25	1.38
2	D	503	3C4	CE2-CD2	-6.57	1.26	1.38
2	A	501	3C4	CE2-CD2	-6.33	1.27	1.38
2	C	502	3C4	CE2-CZ	6.30	1.50	1.39
2	B	504	3C4	CE2-CD2	-5.08	1.29	1.38
2	B	504	3C4	CE2-CZ	4.93	1.48	1.39
2	C	502	3C4	CE2-CD2	-4.58	1.30	1.38
2	B	505	3C4	CD1-CA	-4.39	1.31	1.39
2	D	503	3C4	CD1-CA	-4.28	1.32	1.39
2	A	501	3C4	OH-CZ	4.25	1.45	1.36
2	D	503	3C4	CE2-CZ	3.77	1.46	1.39
2	A	501	3C4	CE2-CZ	3.73	1.46	1.39
2	B	505	3C4	CE2-CZ	3.50	1.45	1.39
2	B	505	3C4	CD1-CE1	-3.24	1.33	1.38
2	B	504	3C4	CD1-CA	-3.06	1.34	1.39
2	B	505	3C4	OH-CZ	2.97	1.42	1.36

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	505	3C4	CE2-CZ-CE1	-9.46	109.37	118.55
2	D	503	3C4	CE2-CZ-CE1	-7.35	111.42	118.55
2	B	504	3C4	CE2-CZ-CE1	-7.06	111.70	118.55
2	C	502	3C4	CE2-CZ-CE1	-6.90	111.85	118.55
2	A	501	3C4	CE2-CZ-CE1	-6.88	111.87	118.55
2	B	505	3C4	CD1-CE1-CZ	5.79	124.27	120.91
2	B	505	3C4	OH-CZ-CE2	3.71	129.43	119.33
2	A	501	3C4	OH-CZ-CE2	2.94	127.32	119.33
2	B	505	3C4	CE2-CD2-CA	2.90	125.02	121.03
2	C	502	3C4	OH-CZ-CE2	2.77	126.85	119.33
2	A	501	3C4	CD2-CE2-CZ	2.66	123.24	120.50
2	A	501	3C4	CE1-CD1-CA	-2.59	118.69	120.46
2	B	504	3C4	CD2-CA-CD1	2.46	121.99	118.54
2	D	503	3C4	OH-CZ-CE2	2.42	125.92	119.33
2	A	501	3C4	CD2-CA-CD1	2.41	121.92	118.54
2	B	505	3C4	CZ-CE1-CL	-2.29	115.22	119.53
2	D	503	3C4	CD2-CA-CD1	2.24	121.68	118.54
2	A	501	3C4	CD1-CE1-CL	2.24	122.10	118.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	504	3C4	OH-CZ-CE2	2.19	125.29	119.33
2	B	505	3C4	CE1-CD1-CA	-2.15	118.99	120.46
2	C	502	3C4	CD2-CA-CD1	2.01	121.36	118.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.