Full wwPDB X-ray Structure Validation Report (i)

Oct 14, 2021 – 04:12 PM EDT

PDB ID	:	3CQU
Title	:	Crystal Structure of Akt-1 complexed with substrate peptide and inhibitor
Authors	:	Pandit, J.
Deposited on	:	2008-04-03
Resolution	:	2.20 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 following versions of software and data (see references (i)) were used in the production of this report:

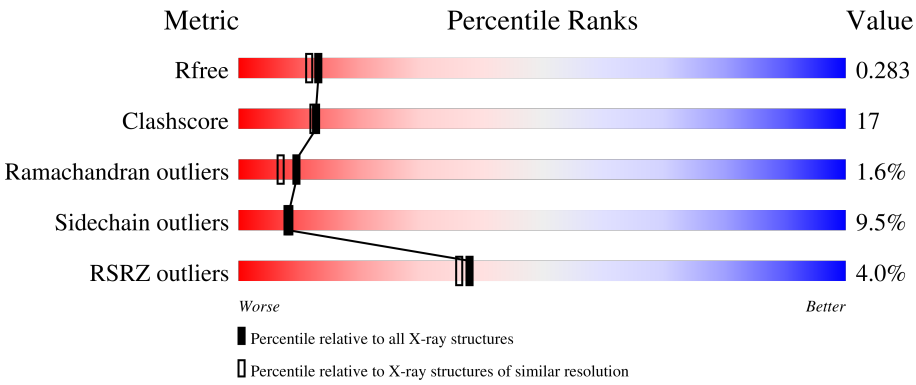
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

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 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	342	4%, 57%, 30%, 7%
2	C	10	10%, 60%, 30%, 10%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2945 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 RAC-alpha serine/threonine-protein kinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	319	2652	1703	447	484	1	17	0	10	0

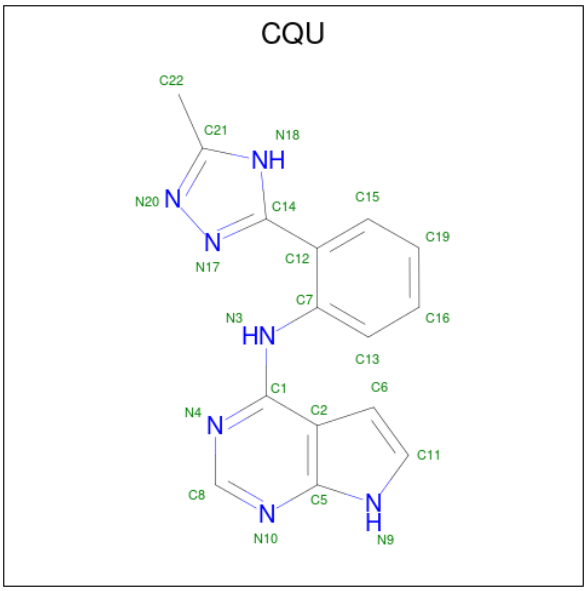
There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	139	GLY	-	expression tag	UNP P31749
A	140	ALA	-	expression tag	UNP P31749
A	141	MET	-	expression tag	UNP P31749
A	142	ASP	-	expression tag	UNP P31749
A	143	PRO	-	expression tag	UNP P31749
A	473	ASP	SER	engineered mutation	UNP P31749
A	478	SER	GLY	variant	UNP P31749

- Molecule 2 is a protein called Glycogen synthase kinase-3 beta.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	10	79	47	16	16	0	0	0

- Molecule 3 is N-[2-(5-methyl-4H-1,2,4-triazol-3-yl)phenyl]-7H-pyrrolo[2,3-d]pyrimidin-4-amine (three-letter code: CQU) (formula: $C_{15}H_{13}N_7$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N 22 15 7	0	0

- Molecule 4 is water.

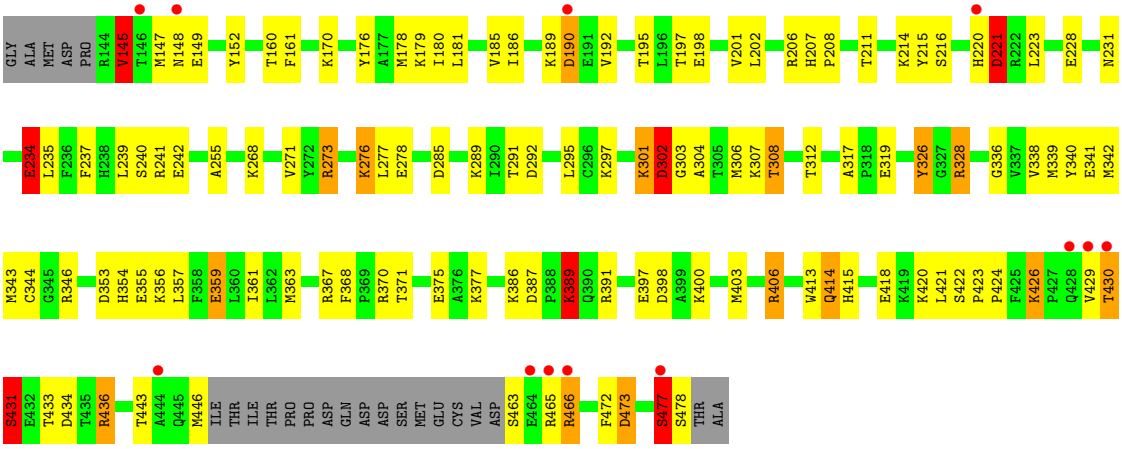
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	184	Total O 184 184	0	0
4	C	8	Total O 8 8	0	0

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: RAC-alpha serine/threonine-protein kinase

Chain A:



• Molecule 2: Glycogen synthase kinase-3 beta

Chain C:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	42.49Å 55.51Å 93.36Å 90.00° 105.24° 90.00°	Depositor
Resolution (Å)	45.04 – 2.20 45.04 – 2.02	Depositor EDS
% Data completeness (in resolution range)	100.0 (45.04-2.20) 99.8 (45.04-2.02)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.93 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.203 , 0.283 0.211 , 0.283	Depositor DCC
R_{free} test set	1392 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	26.9	Xtriage
Anisotropy	0.167	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 53.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.010 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	2945	wwPDB-VP
Average B, all atoms (Å^2)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.66% 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 (i)

5.1 Standard geometry (i)

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

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	1.32	17/2742 (0.6%)	1.20	20/3680 (0.5%)
2	C	1.27	1/80 (1.2%)	1.23	0/105
All	All	1.32	18/2822 (0.6%)	1.20	20/3785 (0.5%)

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	397	GLU	CD-OE2	8.23	1.34	1.25
1	A	341	GLU	CD-OE1	7.60	1.34	1.25
1	A	228	GLU	CD-OE1	7.28	1.33	1.25
1	A	234	GLU	CD-OE1	6.49	1.32	1.25
1	A	397	GLU	CG-CD	6.31	1.61	1.51
1	A	326	TYR	CD2-CE2	-6.17	1.30	1.39
1	A	228	GLU	CD-OE2	6.12	1.32	1.25
1	A	341	GLU	CD-OE2	6.10	1.32	1.25
1	A	391	ARG	CZ-NH1	6.01	1.40	1.33
1	A	278	GLU	CD-OE2	-5.91	1.19	1.25
1	A	336	GLY	C-O	-5.68	1.14	1.23
1	A	389	LYS	CD-CE	5.66	1.65	1.51
1	A	255	ALA	CA-CB	5.37	1.63	1.52
1	A	368	PHE	CE2-CZ	5.35	1.47	1.37
2	C	1	GLY	N-CA	5.18	1.53	1.46
1	A	319	GLU	CD-OE2	-5.15	1.20	1.25
1	A	319	GLU	CD-OE1	5.13	1.31	1.25
1	A	317	ALA	CA-CB	-5.01	1.42	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	190	ASP	CB-CG-OD2	7.80	125.32	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	406	ARG	NE-CZ-NH2	-7.36	116.62	120.30
1	A	473	ASP	CB-CG-OD2	7.13	124.72	118.30
1	A	370	ARG	NE-CZ-NH1	7.04	123.82	120.30
1	A	346	ARG	NE-CZ-NH1	6.59	123.59	120.30
1	A	387	ASP	CB-CG-OD1	6.58	124.23	118.30
1	A	328	ARG	NE-CZ-NH2	-6.40	117.10	120.30
1	A	221	ASP	CB-CG-OD2	6.17	123.85	118.30
1	A	436	ARG	NE-CZ-NH1	6.06	123.33	120.30
1	A	406	ARG	NE-CZ-NH1	5.97	123.28	120.30
1	A	145	VAL	N-CA-C	5.88	126.86	111.00
1	A	302	ASP	CB-CG-OD2	5.86	123.58	118.30
1	A	391	ARG	NE-CZ-NH1	5.77	123.19	120.30
1	A	367	ARG	NE-CZ-NH1	5.73	123.16	120.30
1	A	434	ASP	CB-CG-OD2	5.53	123.28	118.30
1	A	285	ASP	CB-CG-OD2	5.52	123.27	118.30
1	A	273	ARG	NE-CZ-NH2	-5.23	117.69	120.30
1	A	398	ASP	CB-CG-OD2	5.22	123.00	118.30
1	A	190	ASP	OD1-CG-OD2	-5.07	113.67	123.30
1	A	202	LEU	CA-CB-CG	5.01	126.82	115.30

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	2652	0	2618	87	0
2	C	79	0	77	2	0
3	A	22	0	13	4	0
4	A	184	0	0	29	0
4	C	8	0	0	1	0
All	All	2945	0	2708	91	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 17.

All (91) 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:375:GLU:HA	4:A:1176:HOH:O	1.49	1.11
1:A:414[B]:GLN:HG3	1:A:418:GLU:OE2	1.72	0.89
2:C:6:THR:HG22	4:C:19:HOH:O	1.76	0.83
1:A:178:MET:HE3	1:A:180:ILE:HG12	1.62	0.82
1:A:220:HIS:CE1	1:A:221:ASP:HB3	2.15	0.80
1:A:271:VAL:HG21	4:A:1178:HOH:O	1.85	0.77
1:A:328:ARG:HA	4:A:1178:HOH:O	1.84	0.77
1:A:406:ARG:NE	4:A:1176:HOH:O	2.19	0.75
1:A:353:ASP:OD2	1:A:356[B]:LYS:HG2	1.90	0.71
1:A:301:LYS:O	1:A:302:ASP:O	2.08	0.71
1:A:477:SER:HA	4:A:1138:HOH:O	1.90	0.71
1:A:195[A]:THR:HG21	1:A:223:LEU:HD13	1.71	0.70
1:A:400:LYS:HE3	4:A:1085:HOH:O	1.93	0.69
1:A:306:MET:HE1	4:A:1178:HOH:O	1.92	0.68
1:A:148:ASN:O	1:A:170:LYS:NZ	2.19	0.65
1:A:292:ASP:OD2	3:A:999:CQU:H22	1.97	0.65
1:A:161:PHE:O	1:A:179[A]:LYS:HE2	1.97	0.63
1:A:361:ILE:O	1:A:386:LYS:HE3	1.98	0.62
1:A:192:VAL:HG13	4:A:1091:HOH:O	2.00	0.61
1:A:340:TYR:O	1:A:344:CYS:HB2	2.01	0.60
1:A:421:LEU:HD12	4:A:1134:HOH:O	2.01	0.60
1:A:178:MET:HE2	1:A:180:ILE:HD11	1.85	0.59
1:A:211:THR:OG1	1:A:291[B]:THR:HG23	2.05	0.57
1:A:443:THR:HG22	4:A:1144:HOH:O	2.04	0.57
2:C:2:ARG:HH21	2:C:5:THR:HG23	1.70	0.56
1:A:221:ASP:OD1	1:A:221:ASP:O	2.24	0.56
1:A:355:GLU:CD	1:A:356[B]:LYS:NZ	2.60	0.56
1:A:308:TPO:O	1:A:326:TYR:OH	2.17	0.55
1:A:356[A]:LYS:NZ	4:A:1169:HOH:O	2.23	0.55
1:A:239:LEU:HD23	1:A:277:LEU:HD21	1.89	0.55
1:A:355:GLU:HG2	1:A:356[B]:LYS:HD3	1.88	0.55
1:A:328:ARG:CA	4:A:1178:HOH:O	2.47	0.54
1:A:231:ASN:OD1	1:A:429:VAL:HG21	2.08	0.54
1:A:353:ASP:HB3	1:A:356[B]:LYS:HG2	1.91	0.53
1:A:430:THR:O	1:A:431:SER:HB3	2.09	0.53
1:A:339:MET:O	1:A:343[A]:MET:HG3	2.10	0.51
1:A:356[B]:LYS:HD2	4:A:1169:HOH:O	2.11	0.51
1:A:478:SER:CA	4:A:1061:HOH:O	2.58	0.51
1:A:178:MET:CE	1:A:180:ILE:HG12	2.39	0.50
1:A:430:THR:O	1:A:430:THR:HG23	2.11	0.50
1:A:355:GLU:CD	1:A:356[B]:LYS:HZ2	2.14	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:LYS:HE2	1:A:312:THR:OG1	2.12	0.50
1:A:289:LYS:HE3	4:A:1032:HOH:O	2.12	0.50
1:A:145:VAL:HG21	1:A:215:TYR:CD1	2.46	0.50
1:A:271:VAL:HG11	4:A:1178:HOH:O	2.12	0.49
1:A:426:LYS:HD3	4:A:1177:HOH:O	2.10	0.49
1:A:220:HIS:HB2	1:A:466:ARG:NH1	2.27	0.49
3:A:999:CQU:H13	3:A:999:CQU:N4	2.28	0.49
1:A:161:PHE:HB2	1:A:179[A]:LYS:CE	2.43	0.49
1:A:181:LEU:HB3	1:A:186:ILE:HD11	1.95	0.48
1:A:176:TYR:CE1	1:A:214[A]:LYS:HG3	2.47	0.48
1:A:418:GLU:OE1	1:A:420:LYS:HE3	2.13	0.48
1:A:423:PRO:HA	1:A:424:PRO:HD3	1.72	0.47
1:A:237:PHE:O	1:A:240:SER:HB2	2.15	0.47
1:A:415:HIS:HD2	4:A:1149:HOH:O	1.97	0.47
1:A:354:HIS:HA	1:A:357:LEU:HB3	1.97	0.47
1:A:234:GLU:O	1:A:237:PHE:HB3	2.14	0.47
1:A:306:MET:CE	4:A:1178:HOH:O	2.53	0.46
1:A:185:VAL:O	1:A:189:LYS:HG3	2.15	0.46
1:A:235:LEU:HD21	1:A:342[A]:MET:CE	2.45	0.46
1:A:216:SER:HB3	1:A:472:PHE:CE2	2.51	0.46
1:A:465:ARG:CZ	4:A:1116:HOH:O	2.64	0.45
1:A:403:MET:CE	4:A:1010:HOH:O	2.65	0.45
1:A:356[B]:LYS:CD	4:A:1169:HOH:O	2.64	0.44
1:A:415:HIS:CD2	4:A:1149:HOH:O	2.71	0.44
1:A:145:VAL:HG22	4:A:1146:HOH:O	2.18	0.44
1:A:273:ARG:HD3	1:A:295:LEU:O	2.19	0.43
1:A:186:ILE:CD1	1:A:223:LEU:HD12	2.49	0.43
1:A:178:MET:HE3	1:A:180:ILE:CG1	2.42	0.43
1:A:268:LYS:HD3	1:A:268:LYS:HA	1.81	0.43
1:A:403:MET:HE2	4:A:1010:HOH:O	2.18	0.43
1:A:221:ASP:OD1	1:A:221:ASP:C	2.57	0.43
1:A:292:ASP:OD2	3:A:999:CQU:C22	2.65	0.43
1:A:208:PRO:HD3	4:A:1160:HOH:O	2.19	0.42
1:A:406:ARG:CZ	4:A:1176:HOH:O	2.66	0.42
1:A:297:LYS:HG2	1:A:306:MET:HG2	2.01	0.42
1:A:353:ASP:OD2	1:A:356[B]:LYS:HE2	2.20	0.41
1:A:359:GLU:O	1:A:363:MET:HB2	2.21	0.41
1:A:207:HIS:CG	1:A:208:PRO:HD2	2.56	0.41
1:A:338:VAL:O	1:A:342[B]:MET:HG3	2.20	0.41
1:A:152:TYR:CD1	1:A:152:TYR:N	2.87	0.41
1:A:179[B]:LYS:NZ	1:A:198:GLU:OE2	2.40	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:THR:O	1:A:201:VAL:HG23	2.20	0.41
1:A:297:LYS:HD2	1:A:297:LYS:HA	1.89	0.41
1:A:377:LYS:HD3	4:A:1110:HOH:O	2.20	0.41
1:A:178:MET:CE	1:A:180:ILE:HD11	2.51	0.41
1:A:389:LYS:HA	1:A:389:LYS:HD3	1.93	0.40
3:A:999:CQU:H15	4:A:1082:HOH:O	2.21	0.40
1:A:152:TYR:HE2	1:A:178:MET:HE1	1.85	0.40
1:A:403:MET:HA	1:A:413:TRP:CZ2	2.56	0.40
1:A:303:GLY:O	1:A:304:ALA:C	2.59	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	324/342 (95%)	301 (93%)	18 (6%)	5 (2%)	10 8
2	C	8/10 (80%)	8 (100%)	0	0	100 100
All	All	332/352 (94%)	309 (93%)	18 (5%)	5 (2%)	9 8

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	302	ASP
1	A	431	SER
1	A	145	VAL
1	A	477	SER
1	A	190	ASP

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	287/299 (96%)	260 (91%)	27 (9%)	8 8
2	C	8/8 (100%)	7 (88%)	1 (12%)	4 4
All	All	295/307 (96%)	267 (90%)	28 (10%)	8 8

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	147	MET
1	A	149	GLU
1	A	160	THR
1	A	206	ARG
1	A	221	ASP
1	A	234	GLU
1	A	241	ARG
1	A	242	GLU
1	A	276	LYS
1	A	301	LYS
1	A	307	LYS
1	A	359	GLU
1	A	371	THR
1	A	389	LYS
1	A	414[A]	GLN
1	A	414[B]	GLN
1	A	422	SER
1	A	426	LYS
1	A	430	THR
1	A	431	SER
1	A	433	THR
1	A	436	ARG
1	A	446	MET
1	A	463	SER
1	A	466	ARG
1	A	473	ASP
1	A	477	SER

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Mol	Chain	Res	Type
2	C	6	THR

Sometimes 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)

1 non-standard protein/DNA/RNA residue is 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	$\#\|Z\| > 2$
1	TPO	A	308	1	8,10,11	1.62	2 (25%)	10,14,16	0.92	1 (10%)

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
1	TPO	A	308	1	-	1/9/11/13	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	308	TPO	P-O1P	3.06	1.60	1.50
1	A	308	TPO	O-C	2.24	1.28	1.19

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	308	TPO	O3P-P-O2P	2.39	116.77	107.64

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	308	TPO	CB-OG1-P-O2P

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	308	TPO	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand is 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	$\#\|Z\| > 2$
3	CQU	A	999	-	22,25,25	1.15	3 (13%)	24,35,35	3.22	9 (37%)

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
3	CQU	A	999	-	-	1/8/8/8	0/4/4/4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	999	CQU	N17-N20	-3.20	1.31	1.37
3	A	999	CQU	C1-N3	2.81	1.40	1.36
3	A	999	CQU	C5-N10	-2.71	1.33	1.37

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	999	CQU	C14-N17-N20	8.75	112.07	104.70
3	A	999	CQU	C2-C1-N4	-7.03	115.56	121.35
3	A	999	CQU	C8-N4-C1	5.79	121.56	116.59
3	A	999	CQU	C21-N18-C14	5.36	107.15	103.79
3	A	999	CQU	N3-C1-N4	4.41	124.63	118.72
3	A	999	CQU	C12-C14-N17	2.66	128.63	124.11
3	A	999	CQU	C8-N10-C5	2.60	119.53	113.45
3	A	999	CQU	N10-C8-N4	-2.55	124.70	128.68
3	A	999	CQU	C12-C14-N18	2.05	126.29	123.67

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	999	CQU	C7-C12-C14-N17

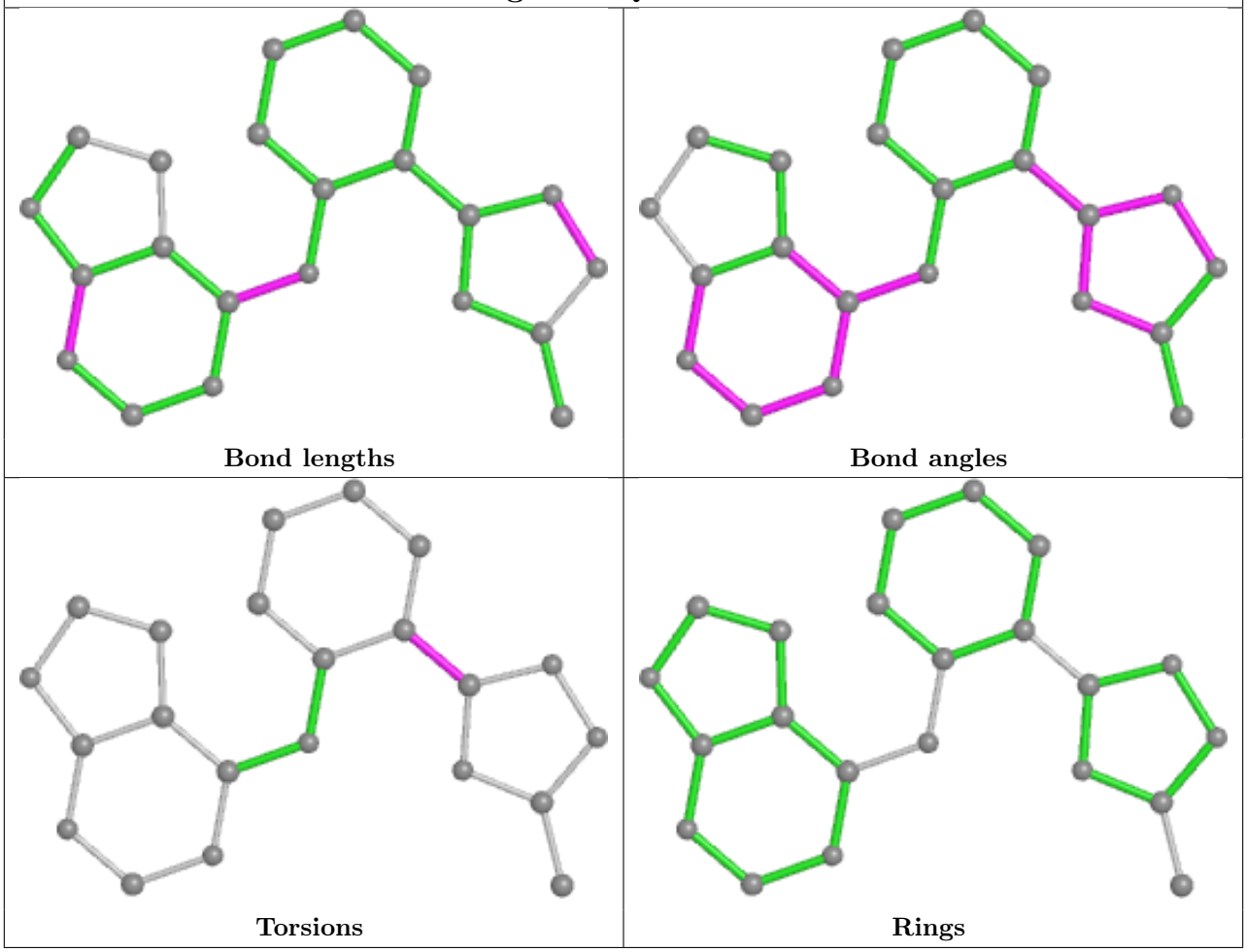
There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	999	CQU	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 CQU A 999



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	318/342 (92%)	-0.11	12 (3%)	40	38	28, 45, 71, 89	0
2	C	10/10 (100%)	0.24	1 (10%)	7	6	45, 51, 56, 62	0
All	All	328/352 (93%)	-0.10	13 (3%)	38	36	28, 45, 71, 89	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	444	ALA	3.9
1	A	430	THR	3.6
2	C	1	GLY	3.5
1	A	477	SER	3.3
1	A	465	ARG	3.0
1	A	148	ASN	2.9
1	A	429	VAL	2.8
1	A	190	ASP	2.7
1	A	464	GLU	2.4
1	A	146	THR	2.3
1	A	466	ARG	2.2
1	A	428	GLN	2.1
1	A	220	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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
1	TPO	A	308	11/12	0.97	0.09	35,40,45,51	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

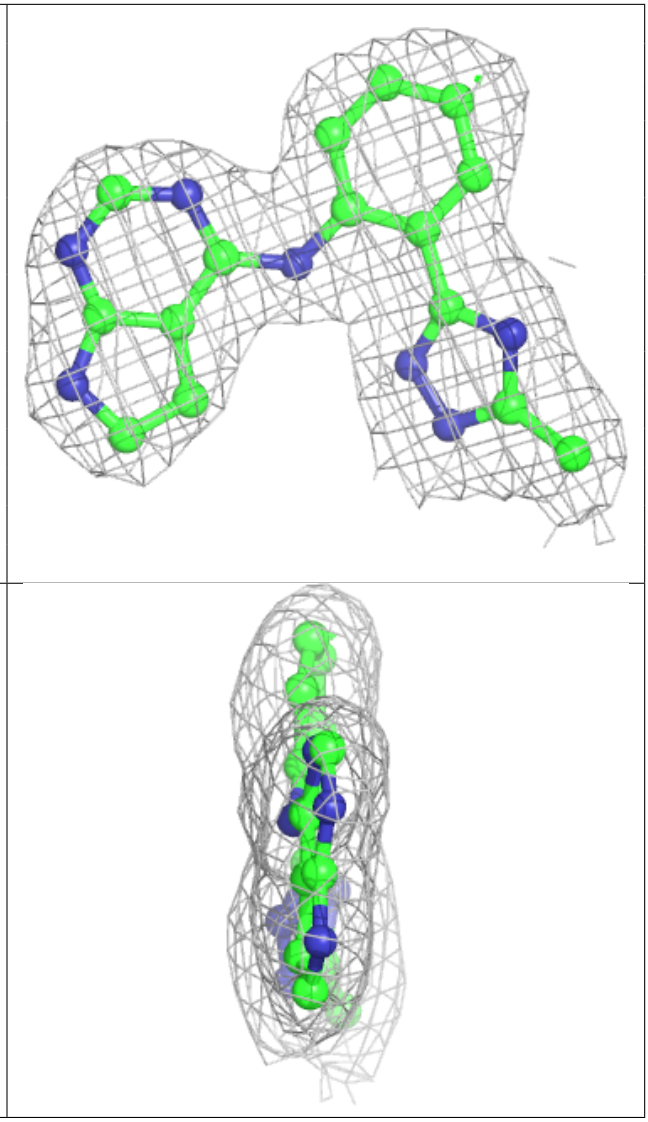
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	CQU	A	999	22/22	0.92	0.13	44,48,51,53	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 CQU A 999:

$2mF_o$-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.