Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 09:07 PM GMT

PDB ID	:	4UYN
Title	:	SAR156497 an exquisitely selective inhibitor of Aurora kinases
Authors	:	Carry, J.C.; Clerc, F.; Minoux, H.; Schio, L.; Mauger, J.; Nair, A.; Parmantier, E.; Lemoigne, R.; Delorme, C.; Nicolas, J.P.; Krick, A.; Abecassis, P.Y.; Crocq-Stuerga, V.; Pouzieux, S.; Delarbre, L.; Maignan, S.; Bertrand, T.; Bjergarde, K.; Ma, N.; Lachaud, S.; Guizani, H.; Lebel, R.; Doerflinger, G.; Monget, S.; Perron, S.; Gasse, F.; Angouillant-Boniface, O.; Filoche-Romme, B.; Murer, M.; Gontier, S.; Prevost, C.; Monteiro, M.L.; Combeau, C.
Deposited on	:	2014-09-02
Resolution	:	1.90 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

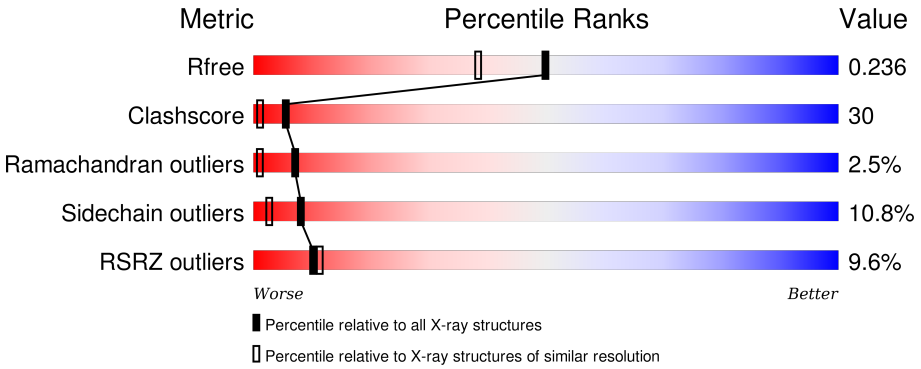
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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 1.90 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}	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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. 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	287	8% 53% 26% 6% • 13%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2182 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.

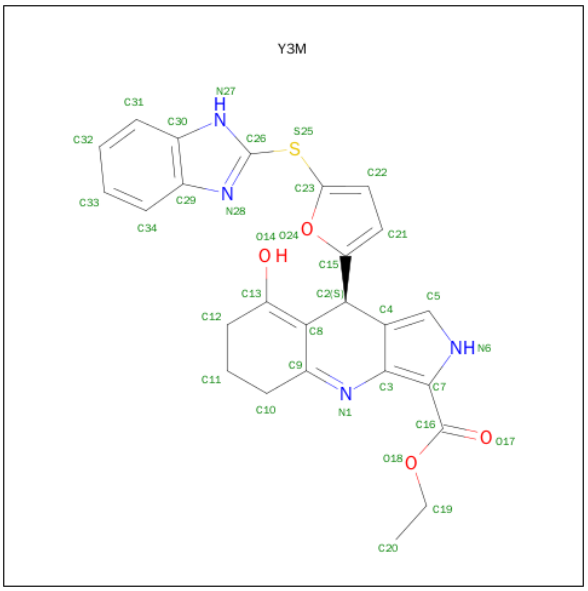
- Molecule 1 is a protein called AURORA KINASE A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	0	0	0
			2059	1327	361	366	5			

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	124	MET	-	EXPRESSION TAG	UNP O14965
A	400	ALA	-	EXPRESSION TAG	UNP O14965
A	401	ALA	-	EXPRESSION TAG	UNP O14965
A	402	ALA	-	EXPRESSION TAG	UNP O14965
A	403	LEU	-	EXPRESSION TAG	UNP O14965
A	404	GLU	-	EXPRESSION TAG	UNP O14965
A	405	HIS	-	EXPRESSION TAG	UNP O14965
A	406	HIS	-	EXPRESSION TAG	UNP O14965
A	407	HIS	-	EXPRESSION TAG	UNP O14965
A	408	HIS	-	EXPRESSION TAG	UNP O14965
A	409	HIS	-	EXPRESSION TAG	UNP O14965
A	410	HIS	-	EXPRESSION TAG	UNP O14965
A	288	ASP	THR	CONFLICT	UNP O14965

- Molecule 2 is ETHYL (9S)-9-[5-(1H-BENZIMIDAZOL-2-YLSULFANYL)FURAN-2-YL]-8-HYDROXY-5,6,7,9-TETRAHYDRO-2H-PYRROLO[3,4-B]QUINOLINE-3-CARBOXYLATE (three-letter code: Y3M) (formula: $C_{25}H_{22}N_4O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			34	25	4	4	1		

- Molecule 3 is water.

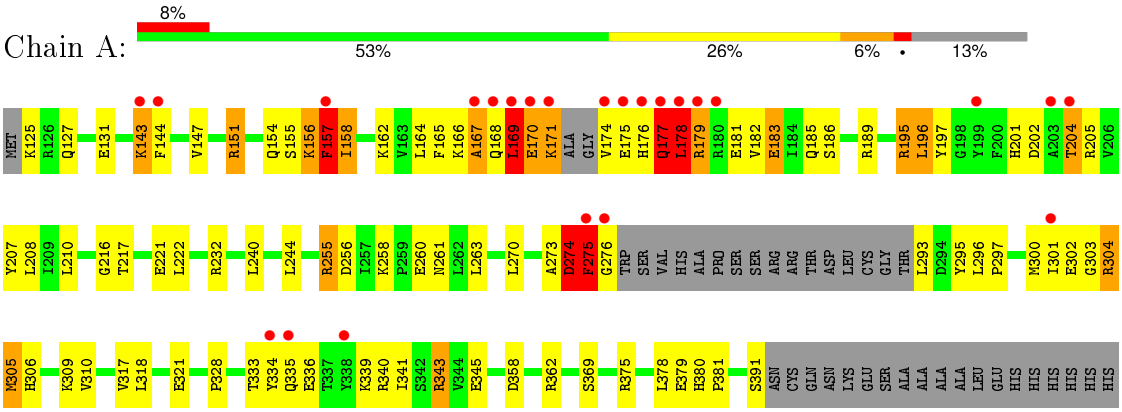
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	89	Total	O	0	0
			89	89		

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 errors 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: AURORA KINASE A

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	48.35Å 45.91Å 59.59Å 90.00° 97.82° 90.00°	Depositor
Resolution (Å)	59.00 – 1.90 59.04 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.2 (59.00-1.90) 99.2 (59.04-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.32 (at 1.90Å)	Xtriage
Refinement program	BUSTER-TNT 2.9.3	Depositor
R, R_{free}	0.201 , 0.234 0.207 , 0.236	Depositor DCC
R_{free} test set	1607 reflections (8.52%)	DCC
Wilson B-factor (Å^2)	22.7	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 66.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 20466 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	2182	wwPDB-VP
Average B, all atoms (Å^2)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.01% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: Y3M

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	0.78	4/2108 (0.2%)	0.81	8/2843 (0.3%)

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	5	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	304	ARG	C-N	-20.86	0.86	1.34
1	A	157	PHE	C-N	-13.06	1.04	1.34
1	A	303	GLY	C-N	12.26	1.62	1.34
1	A	158	ILE	C-N	5.71	1.47	1.34

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	303	GLY	O-C-N	-10.36	106.13	122.70
1	A	156	LYS	O-C-N	8.99	137.09	122.70
1	A	156	LYS	C-N-CA	-7.65	102.58	121.70
1	A	303	GLY	CA-C-N	7.10	132.81	117.20
1	A	156	LYS	CA-C-N	-7.07	101.65	117.20
1	A	178	LEU	N-CA-C	6.68	129.04	111.00
1	A	303	GLY	C-N-CA	6.57	138.12	121.70
1	A	157	PHE	CA-C-N	-5.56	104.96	117.20

All (5) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	125	LYS	CA
1	A	174	VAL	CA
1	A	183	GLU	CA
1	A	274	ASP	CA
1	A	275	PHE	CA

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	157	PHE	Mainchain

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	2059	0	2072	124	1
2	A	34	0	21	8	0
3	A	89	0	0	8	0
All	All	2182	0	2093	124	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 30.

All (124) 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:304:ARG:C	1:A:305:MET:CA	1.98	1.31
1:A:304:ARG:CA	1:A:305:MET:N	1.96	1.28
1:A:304:ARG:O	1:A:305:MET:N	1.91	1.02
1:A:176:HIS:C	1:A:179:ARG:HB2	1.85	0.95
1:A:300:MET:HE2	1:A:310:VAL:HG11	1.47	0.93
1:A:168:GLN:HA	1:A:171:LYS:HG3	1.52	0.91
1:A:304:ARG:C	1:A:305:MET:N	0.86	0.90
1:A:302:GLU:OE1	1:A:304:ARG:NH1	2.09	0.85
1:A:273:ALA:O	1:A:274:ASP:O	1.93	0.84

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:GLN:O	1:A:177:GLN:HG2	1.77	0.84
1:A:167:ALA:O	1:A:170:GLU:HG3	1.83	0.78
1:A:300:MET:CE	1:A:310:VAL:HG11	2.15	0.77
1:A:339:LYS:HG3	1:A:343:ARG:HD2	1.68	0.76
1:A:169:LEU:O	1:A:174:VAL:HB	1.86	0.76
1:A:178:LEU:HD21	1:A:179:ARG:NH1	2.02	0.75
1:A:221:GLU:HG2	1:A:232:ARG:NH1	2.00	0.75
1:A:143:LYS:HD2	1:A:143:LYS:O	1.87	0.74
1:A:151:ARG:NH1	1:A:156:LYS:HD2	2.02	0.73
1:A:183:GLU:HB2	1:A:186:SER:OG	1.88	0.73
1:A:306:HIS:HD2	3:A:2056:HOH:O	1.71	0.72
1:A:258:LYS:HE2	1:A:261:ASN:ND2	2.05	0.72
1:A:177:GLN:N	1:A:179:ARG:H	1.88	0.72
1:A:178:LEU:HD23	1:A:179:ARG:N	2.06	0.71
1:A:178:LEU:HD23	1:A:179:ARG:H	1.55	0.70
1:A:258:LYS:HE2	1:A:261:ASN:HD21	1.56	0.70
1:A:168:GLN:CA	1:A:171:LYS:HG3	2.22	0.69
1:A:168:GLN:HB3	1:A:171:LYS:HG3	1.72	0.69
1:A:168:GLN:CB	1:A:171:LYS:HG3	2.23	0.68
1:A:181:GLU:HG3	1:A:276:GLY:HA2	1.74	0.68
1:A:183:GLU:HA	1:A:183:GLU:OE1	1.93	0.68
1:A:195:ARG:HG3	1:A:195:ARG:NH1	2.09	0.68
1:A:304:ARG:O	1:A:305:MET:CA	2.34	0.68
1:A:205:ARG:HD2	1:A:207:TYR:OH	1.98	0.64
1:A:256:ASP:OD1	1:A:258:LYS:HG2	1.96	0.64
1:A:216:GLY:HA2	2:A:1392:Y3M:H201	1.78	0.64
1:A:178:LEU:HD21	1:A:179:ARG:HH11	1.61	0.64
1:A:176:HIS:O	1:A:177:GLN:HB3	1.96	0.64
1:A:375:ARG:O	1:A:379:GLU:HG3	2.00	0.62
1:A:275:PHE:HB2	2:A:1392:Y3M:N28	2.17	0.59
1:A:155:SER:O	1:A:156:LYS:HB2	2.02	0.59
1:A:358:ASP:O	1:A:362:ARG:HG3	2.02	0.59
1:A:263:LEU:HD22	2:A:1392:Y3M:C20	2.33	0.59
1:A:176:HIS:HA	1:A:179:ARG:CG	2.32	0.59
1:A:165:PHE:HA	1:A:204:THR:O	2.02	0.59
1:A:157:PHE:CG	1:A:158:ILE:N	2.69	0.59
1:A:176:HIS:CA	1:A:179:ARG:HB2	2.32	0.59
1:A:195:ARG:HH11	1:A:195:ARG:HG3	1.67	0.58
1:A:263:LEU:HD22	2:A:1392:Y3M:H203	1.86	0.57
1:A:293:LEU:N	3:A:2049:HOH:O	2.37	0.57
1:A:170:GLU:C	1:A:171:LYS:HG2	2.25	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:GLN:O	1:A:177:GLN:CG	2.51	0.56
1:A:195:ARG:HG2	1:A:197:TYR:CE1	2.41	0.56
1:A:343:ARG:NH2	3:A:2064:HOH:O	2.39	0.55
1:A:127:GLN:NE2	3:A:2003:HOH:O	2.40	0.55
1:A:336:GLU:O	1:A:340:ARG:HG2	2.07	0.55
1:A:240:LEU:CD1	1:A:318:LEU:HD23	2.37	0.54
1:A:258:LYS:CE	1:A:261:ASN:HD21	2.18	0.54
1:A:296:LEU:HD23	3:A:2042:HOH:O	2.06	0.54
1:A:302:GLU:OE1	1:A:304:ARG:NH2	2.41	0.53
1:A:168:GLN:HA	1:A:171:LYS:CG	2.32	0.52
1:A:240:LEU:HD13	1:A:318:LEU:HD23	1.92	0.52
1:A:297:PRO:HG2	1:A:300:MET:HG3	1.92	0.51
1:A:157:PHE:CE2	1:A:158:ILE:O	2.63	0.51
1:A:343:ARG:NH1	1:A:345:GLU:OE1	2.44	0.51
1:A:273:ALA:O	1:A:274:ASP:C	2.49	0.51
1:A:196:LEU:HD12	1:A:210:LEU:HD21	1.92	0.51
1:A:196:LEU:HD12	1:A:210:LEU:CD2	2.40	0.51
1:A:244:LEU:CD2	1:A:255:ARG:HG3	2.41	0.51
1:A:176:HIS:HA	1:A:179:ARG:HG2	1.91	0.50
1:A:302:GLU:OE1	1:A:304:ARG:CZ	2.59	0.50
1:A:177:GLN:HA	1:A:179:ARG:N	2.26	0.49
1:A:309:LYS:NZ	3:A:2057:HOH:O	2.28	0.49
1:A:177:GLN:H	1:A:178:LEU:HD23	1.77	0.49
1:A:151:ARG:HH11	1:A:156:LYS:HD2	1.77	0.49
1:A:177:GLN:N	1:A:179:ARG:HB2	2.27	0.48
1:A:304:ARG:CB	1:A:305:MET:N	2.71	0.48
1:A:165:PHE:O	1:A:169:LEU:HB2	2.13	0.48
1:A:240:LEU:CD1	1:A:318:LEU:CD2	2.92	0.48
1:A:300:MET:HE2	1:A:310:VAL:CG1	2.32	0.48
1:A:258:LYS:HG3	1:A:260:GLU:HG2	1.96	0.48
1:A:177:GLN:CA	1:A:179:ARG:H	2.26	0.47
1:A:178:LEU:O	1:A:181:GLU:HB2	2.14	0.47
1:A:166:LYS:NZ	1:A:201:HIS:HD2	2.13	0.47
1:A:221:GLU:CG	1:A:232:ARG:NH1	2.75	0.47
1:A:317:VAL:HG13	1:A:328:PRO:HD2	1.96	0.46
1:A:178:LEU:HD21	1:A:179:ARG:HD3	1.98	0.46
1:A:221:GLU:HG2	1:A:232:ARG:HH11	1.77	0.46
1:A:166:LYS:HG3	1:A:204:THR:C	2.37	0.45
1:A:306:HIS:HB2	1:A:310:VAL:CG1	2.47	0.45
1:A:166:LYS:HG3	1:A:204:THR:CA	2.46	0.45
1:A:178:LEU:CD2	1:A:179:ARG:HD3	2.47	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:HIS:HA	1:A:179:ARG:HB2	1.97	0.45
1:A:275:PHE:O	1:A:276:GLY:C	2.55	0.45
1:A:195:ARG:HH11	1:A:195:ARG:CG	2.27	0.44
1:A:183:GLU:O	1:A:186:SER:OG	2.31	0.44
1:A:177:GLN:CA	1:A:179:ARG:N	2.81	0.44
1:A:178:LEU:CD2	1:A:179:ARG:N	2.78	0.44
1:A:301:ILE:HD11	1:A:341:ILE:HG22	1.99	0.44
1:A:182:VAL:HG13	1:A:196:LEU:HD21	2.00	0.44
1:A:144:PHE:HB2	1:A:164:LEU:HD22	1.99	0.43
1:A:306:HIS:CD2	3:A:2056:HOH:O	2.58	0.42
1:A:166:LYS:NZ	1:A:202:ASP:O	2.37	0.42
1:A:221:GLU:HG2	1:A:232:ARG:HH12	1.80	0.42
1:A:144:PHE:CD2	1:A:169:LEU:CD2	3.02	0.42
1:A:240:LEU:HG	1:A:270:LEU:HD22	2.02	0.42
1:A:304:ARG:C	1:A:305:MET:C	2.70	0.42
1:A:155:SER:O	1:A:156:LYS:CB	2.68	0.42
1:A:240:LEU:HD11	1:A:318:LEU:CD2	2.50	0.41
1:A:301:ILE:HD13	1:A:341:ILE:HG21	2.03	0.41
1:A:295:TYR:OH	1:A:321:GLU:OE1	2.22	0.41
1:A:263:LEU:HD22	2:A:1392:Y3M:H202	2.03	0.41
1:A:164:LEU:HD23	3:A:2010:HOH:O	2.20	0.41
1:A:162:LYS:HZ3	2:A:1392:Y3M:H27	1.67	0.41
1:A:333:THR:OG1	1:A:336:GLU:HG3	2.21	0.41
1:A:378:LEU:HD22	1:A:391:SER:HB2	2.02	0.41
1:A:170:GLU:OE1	1:A:171:LYS:HE2	2.21	0.41
1:A:334:TYR:CG	1:A:335:GLN:N	2.89	0.41
1:A:147:VAL:HG21	2:A:1392:Y3M:C13	2.50	0.41
1:A:178:LEU:HA	1:A:181:GLU:HB2	2.03	0.41
1:A:162:LYS:HB3	1:A:208:LEU:HB2	2.04	0.40
1:A:240:LEU:HD11	1:A:318:LEU:HD21	2.03	0.40
1:A:185:GLN:OE1	2:A:1392:Y3M:N28	2.54	0.40
1:A:217:THR:HG22	1:A:263:LEU:HD23	2.03	0.40
1:A:380:HIS:CG	1:A:381:PRO:HD2	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:PHE:CZ	1:A:334:TYR:OH[1_545]	2.03	0.17

5.3 Torsion angles

5.3.1 Protein backbone

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	243/287 (85%)	224 (92%)	13 (5%)	6 (2%)	7 1

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	167	ALA
1	A	169	LEU
1	A	177	GLN
1	A	178	LEU
1	A	274	ASP
1	A	275	PHE

5.3.2 Protein sidechains

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	222/252 (88%)	198 (89%)	24 (11%)	8 3

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	125	LYS
1	A	131	GLU
1	A	143	LYS
1	A	151	ARG
1	A	154	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	169	LEU
1	A	170	GLU
1	A	171	LYS
1	A	175	GLU
1	A	177	GLN
1	A	178	LEU
1	A	179	ARG
1	A	183	GLU
1	A	189	ARG
1	A	195	ARG
1	A	196	LEU
1	A	204	THR
1	A	222	LEU
1	A	255	ARG
1	A	274	ASP
1	A	275	PHE
1	A	305	MET
1	A	343	ARG
1	A	369	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	201	HIS
1	A	254	HIS
1	A	261	ASN
1	A	332	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no carbohydrates 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	$\#\|Z\| > 2$
2	Y3M	A	1392	-	31,39,39	2.36	7 (22%)	31,57,57	2.27	12 (38%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	Y3M	A	1392	-	-	0/7/41/41	0/4/6/6

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1392	Y3M	C4-C2	-5.90	1.45	1.52
2	A	1392	Y3M	O14-C13	-4.20	1.22	1.32
2	A	1392	Y3M	C2-C8	-3.71	1.47	1.52
2	A	1392	Y3M	C5-C4	2.17	1.42	1.38
2	A	1392	Y3M	C8-C9	2.93	1.50	1.44
2	A	1392	Y3M	O18-C16	5.17	1.45	1.33
2	A	1392	Y3M	C8-C13	7.04	1.50	1.37

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1392	Y3M	C3-C7-C16	-6.20	119.15	131.61
2	A	1392	Y3M	C12-C13-C8	-3.58	119.45	122.72
2	A	1392	Y3M	C4-C2-C15	-3.26	106.61	111.46
2	A	1392	Y3M	O18-C16-O17	-2.79	118.54	123.66
2	A	1392	Y3M	O14-C13-C8	-2.36	119.99	123.25
2	A	1392	Y3M	C21-C22-C23	-2.22	104.32	107.27
2	A	1392	Y3M	C5-N6-C7	2.11	109.79	104.34

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1392	Y3M	O18-C16-C7	2.55	119.45	112.49
2	A	1392	Y3M	O14-C13-C12	2.66	120.55	114.45
2	A	1392	Y3M	C13-C8-C9	2.82	119.83	118.03
2	A	1392	Y3M	C4-C2-C8	3.51	110.97	105.47
2	A	1392	Y3M	C16-C7-N6	5.15	130.51	120.29

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1392	Y3M	8	0

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

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, 95^{th} 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	249/287 (86%)	0.49	24 (9%)	10	11	10, 24, 67, 106	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	178	LEU	7.2
1	A	157	PHE	5.8
1	A	168	GLN	5.6
1	A	334	TYR	5.5
1	A	338	TYR	5.5
1	A	204	THR	5.4
1	A	276	GLY	5.3
1	A	169	LEU	4.9
1	A	203	ALA	4.7
1	A	170	GLU	4.3
1	A	179	ARG	4.1
1	A	174	VAL	4.0
1	A	144	PHE	3.6
1	A	171	LYS	3.5
1	A	143	LYS	3.4
1	A	176	HIS	3.4
1	A	167	ALA	3.2
1	A	175	GLU	2.8
1	A	301	ILE	2.7
1	A	180	ARG	2.6
1	A	275	PHE	2.4
1	A	177	GLN	2.3
1	A	335	GLN	2.3
1	A	199	TYR	2.2

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} 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	LLDF	B-factors(Å^2)	Q<0.9
2	Y3M	A	1392	34/34	0.96	0.10	-0.62	15,23,27,31	0

6.5 Other polymers (i)

There are no such residues in this entry.