Full wwPDB X-ray Structure Validation Report

Feb 1, 2016 – 07:25 AM GMT

PDB ID	:	3ARX
Title	:	Crystal Structure Analysis of Chitinase A from Vibrio harveyi with novel inhibitors - complex structure with Propentofylline
Authors	:	Pantoom, S.; Vetter, I.R.; Prinz, H.; Suginta, W.
Deposited on	:	2010-12-09
Resolution	:	1.16 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:

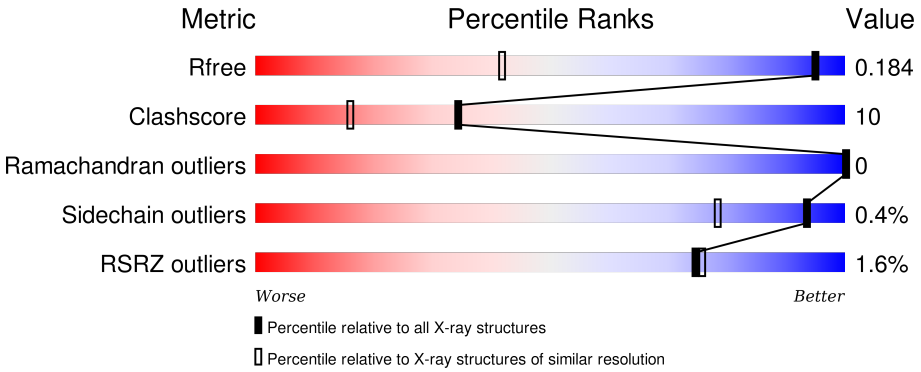
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 1.16 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	1036 (1.20-1.12)
Clashscore	102246	1109 (1.20-1.12)
Ramachandran outliers	100387	1058 (1.20-1.12)
Sidechain outliers	100360	1058 (1.20-1.12)
RSRZ outliers	91569	1038 (1.20-1.12)

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	584	2% 80% 17%

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
2	POY	A	606	-	-	-	X

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	POY	A	607[A]	-	-	X	X
2	POY	A	607[B]	-	-	-	X
3	GOL	A	3500	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5759 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.

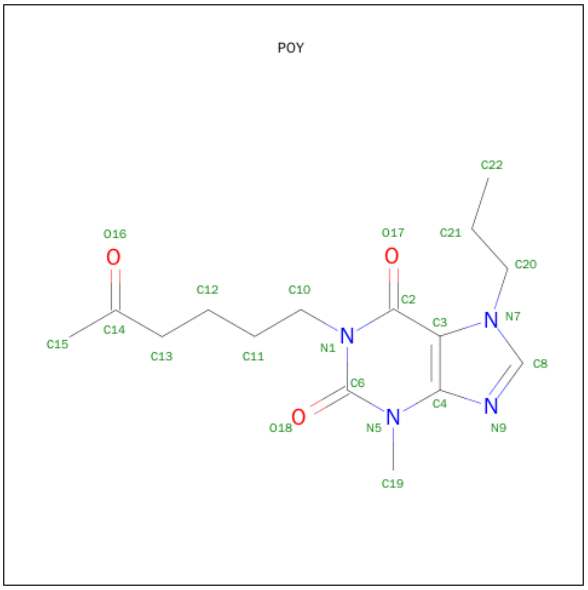
- Molecule 1 is a protein called Chitinase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	575	4588	2931	728	900	29	0	34	0

There are 8 discrepancies between the modelled and reference sequences:

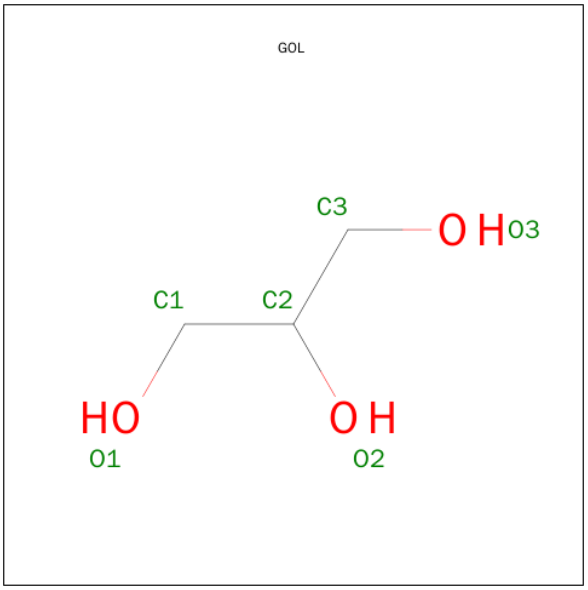
Chain	Residue	Modelled	Actual	Comment	Reference
A	598	ARG	-	EXPRESSION TAG	UNP Q9AMP1
A	599	SER	-	EXPRESSION TAG	UNP Q9AMP1
A	600	HIS	-	EXPRESSION TAG	UNP Q9AMP1
A	601	HIS	-	EXPRESSION TAG	UNP Q9AMP1
A	602	HIS	-	EXPRESSION TAG	UNP Q9AMP1
A	603	HIS	-	EXPRESSION TAG	UNP Q9AMP1
A	604	HIS	-	EXPRESSION TAG	UNP Q9AMP1
A	605	HIS	-	EXPRESSION TAG	UNP Q9AMP1

- Molecule 2 is 3-METHYL-1-(5-OXOHEXYL)-7-PROPYL-3,7-DIHYDRO-1H-PURINE-2,6-DIONE (three-letter code: POY) (formula: $C_{15}H_{22}N_4O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C N O 22 15 4 3	0	0
2	A	1	Total C N O 44 30 8 6	0	1

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0

- Molecule 4 is water.

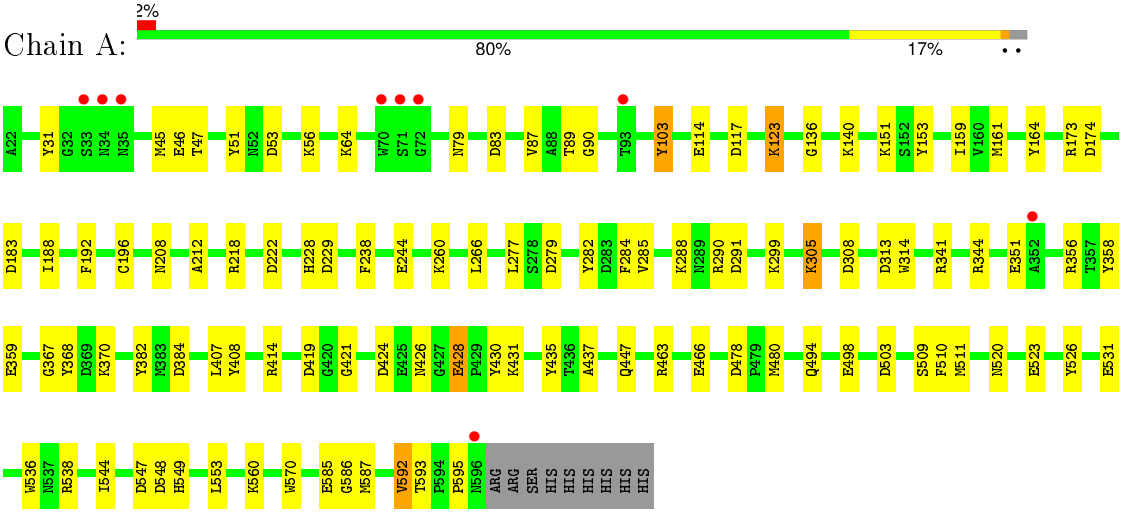
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1099	Total O 1099 1099	0	0

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: Chitinase 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	65.11Å 50.93Å 93.17Å 90.00° 99.45° 90.00°	Depositor
Resolution (Å)	40.66 – 1.16 40.67 – 1.16	Depositor EDS
% Data completeness (in resolution range)	98.4 (40.66-1.16) 98.4 (40.67-1.16)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.11 (at 1.16Å)	Xtriage
Refinement program	REFMAC 5.6.0093	Depositor
R, R_{free}	0.161 , 0.185 0.160 , 0.184	Depositor DCC
R_{free} test set	10234 reflections (5.26%)	DCC
Wilson B-factor (Å^2)	10.6	Xtriage
Anisotropy	0.350	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 45.8	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 204668 reflections	Xtriage
F_o,F_c correlation	0.98	EDS
Total number of atoms	5759	wwPDB-VP
Average B, all atoms (Å^2)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.91% 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: GOL, POY

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	1.64	35/4796 (0.7%)	1.46	60/6525 (0.9%)

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	153	TYR	CE2-CZ	8.84	1.50	1.38
1	A	523	GLU	CB-CG	-8.60	1.35	1.52
1	A	509	SER	CB-OG	-8.34	1.31	1.42
1	A	382	TYR	CE2-CZ	-7.88	1.28	1.38
1	A	466	GLU	CG-CD	7.13	1.62	1.51
1	A	510	PHE	CE1-CZ	-7.02	1.24	1.37
1	A	430	TYR	CD1-CE1	-6.74	1.29	1.39
1	A	123	LYS	CB-CG	-6.66	1.34	1.52
1	A	51	TYR	CE1-CZ	-6.53	1.30	1.38
1	A	164[A]	TYR	CE2-CZ	-6.33	1.30	1.38
1	A	164[B]	TYR	CE2-CZ	-6.33	1.30	1.38
1	A	408	TYR	CE1-CZ	-6.28	1.30	1.38
1	A	368	TYR	CE1-CZ	-6.18	1.30	1.38
1	A	531	GLU	CD-OE2	6.03	1.32	1.25
1	A	408	TYR	CD2-CE2	6.00	1.48	1.39
1	A	103	TYR	CD1-CE1	-5.97	1.30	1.39
1	A	494	GLN	CA-CB	-5.92	1.41	1.53
1	A	153	TYR	CG-CD2	-5.87	1.31	1.39
1	A	359	GLU	CD-OE1	-5.84	1.19	1.25
1	A	196	CYS	CB-SG	5.77	1.92	1.82
1	A	344	ARG	CZ-NH2	5.58	1.40	1.33
1	A	466	GLU	CB-CG	-5.55	1.41	1.52
1	A	592[A]	VAL	CB-CG1	5.51	1.64	1.52
1	A	592[B]	VAL	CB-CG1	5.51	1.64	1.52
1	A	103	TYR	CG-CD1	5.47	1.46	1.39
1	A	382	TYR	CG-CD1	-5.46	1.32	1.39

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	536	TRP	CG-CD1	-5.44	1.29	1.36
1	A	447	GLN	CA-CB	-5.38	1.42	1.53
1	A	430	TYR	CE1-CZ	5.31	1.45	1.38
1	A	560	LYS	CE-NZ	5.31	1.62	1.49
1	A	238	PHE	CG-CD2	-5.26	1.30	1.38
1	A	314	TRP	CG-CD1	-5.26	1.29	1.36
1	A	136	GLY	CA-C	5.23	1.60	1.51
1	A	114	GLU	CD-OE2	-5.17	1.20	1.25
1	A	421	GLY	C-O	5.01	1.31	1.23

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	290	ARG	NE-CZ-NH2	-12.78	113.91	120.30
1	A	218	ARG	NE-CZ-NH2	-12.60	114.00	120.30
1	A	538	ARG	NE-CZ-NH1	-11.62	114.49	120.30
1	A	222	ASP	CB-CG-OD2	-11.59	107.87	118.30
1	A	218	ARG	NE-CZ-NH1	11.57	126.08	120.30
1	A	341[A]	ARG	NE-CZ-NH2	-10.48	115.06	120.30
1	A	341[B]	ARG	NE-CZ-NH2	-10.48	115.06	120.30
1	A	183	ASP	CB-CG-OD2	-9.96	109.34	118.30
1	A	290	ARG	NE-CZ-NH1	9.82	125.21	120.30
1	A	419	ASP	CB-CG-OD2	9.64	126.98	118.30
1	A	51	TYR	CB-CG-CD2	-8.96	115.62	121.00
1	A	173	ARG	NE-CZ-NH1	8.37	124.48	120.30
1	A	222	ASP	CB-CG-OD1	8.26	125.74	118.30
1	A	31	TYR	CB-CG-CD1	-8.02	116.19	121.00
1	A	503	ASP	CB-CG-OD1	7.87	125.38	118.30
1	A	291	ASP	CB-CG-OD1	7.76	125.29	118.30
1	A	548	ASP	CB-CG-OD1	7.58	125.12	118.30
1	A	419	ASP	CB-CG-OD1	-7.41	111.63	118.30
1	A	279	ASP	CB-CG-OD1	7.40	124.96	118.30
1	A	45	MET	N-CA-C	7.25	130.58	111.00
1	A	103	TYR	CB-CG-CD2	7.23	125.34	121.00
1	A	384	ASP	CB-CG-OD2	7.17	124.75	118.30
1	A	585[A]	GLU	OE1-CD-OE2	-7.10	114.78	123.30
1	A	585[B]	GLU	OE1-CD-OE2	-7.10	114.78	123.30
1	A	351	GLU	OE1-CD-OE2	-7.00	114.89	123.30
1	A	463	ARG	NE-CZ-NH1	6.89	123.75	120.30
1	A	31	TYR	CB-CG-CD2	6.83	125.10	121.00
1	A	435	TYR	CB-CG-CD2	-6.78	116.93	121.00
1	A	308	ASP	CB-CG-OD2	6.50	124.15	118.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	83	ASP	CB-CG-OD1	-6.38	112.55	118.30
1	A	284	PHE	CB-CG-CD1	6.37	125.26	120.80
1	A	284	PHE	CB-CG-CD2	-6.36	116.35	120.80
1	A	282	TYR	CB-CG-CD1	-6.29	117.22	121.00
1	A	435	TYR	CB-CG-CD1	6.19	124.71	121.00
1	A	547	ASP	CB-CG-OD1	6.16	123.84	118.30
1	A	358	TYR	CB-CG-CD2	6.11	124.67	121.00
1	A	356	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	A	53[A]	ASP	CB-CG-OD2	-5.94	112.96	118.30
1	A	53[B]	ASP	CB-CG-OD2	-5.94	112.96	118.30
1	A	51	TYR	CG-CD2-CE2	-5.92	116.56	121.30
1	A	358	TYR	CG-CD1-CE1	5.82	125.96	121.30
1	A	53[A]	ASP	CB-CG-OD1	5.76	123.48	118.30
1	A	53[B]	ASP	CB-CG-OD1	5.76	123.48	118.30
1	A	117	ASP	CB-CG-OD1	5.75	123.47	118.30
1	A	159[A]	ILE	CA-CB-CG2	5.66	122.22	110.90
1	A	159[B]	ILE	CA-CB-CG2	5.66	122.22	110.90
1	A	279	ASP	CB-CG-OD2	-5.64	113.22	118.30
1	A	478	ASP	CB-CG-OD2	-5.59	113.27	118.30
1	A	480[A]	MET	CG-SD-CE	5.57	109.10	100.20
1	A	480[B]	MET	CG-SD-CE	5.57	109.10	100.20
1	A	341[A]	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	A	341[B]	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	A	282	TYR	CB-CG-CD2	5.50	124.30	121.00
1	A	260	LYS	CD-CE-NZ	5.41	124.14	111.70
1	A	313	ASP	CB-CG-OD1	5.35	123.11	118.30
1	A	526	TYR	CB-CG-CD1	-5.14	117.91	121.00
1	A	103	TYR	CB-CG-CD1	-5.13	117.92	121.00
1	A	183	ASP	CB-CG-OD1	5.11	122.89	118.30
1	A	305	LYS	CD-CE-NZ	-5.02	100.16	111.70
1	A	114	GLU	N-CA-CB	-5.01	101.59	110.60

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	4588	0	4406	74	2
2	A	66	0	66	28	0
3	A	6	0	8	1	0
4	A	1099	0	0	35	7
All	All	5759	0	4480	91	7

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 10.

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:305:LYS:HD2	4:A:904:HOH:O	1.28	1.30
2:A:607[A]:POY:H15	4:A:1276:HOH:O	1.27	1.26
2:A:607[A]:POY:H10	4:A:949:HOH:O	1.43	1.19
1:A:151:LYS:HE2	4:A:1236:HOH:O	1.41	1.19
1:A:498:GLU:OE1	2:A:607[B]:POY:H15B	1.44	1.15
1:A:229:ASP:HA	2:A:607[A]:POY:H15B	1.33	1.08
2:A:607[B]:POY:O17	2:A:607[B]:POY:H20	1.53	1.05
2:A:607[A]:POY:H8	2:A:607[A]:POY:H22A	1.39	1.04
1:A:511[B]:MET:SD	1:A:544:ILE:CD1	2.48	1.01
1:A:511[B]:MET:SD	1:A:544:ILE:HD12	2.02	1.00
2:A:607[A]:POY:C15	4:A:1276:HOH:O	1.89	1.00
2:A:607[A]:POY:O16	4:A:1595:HOH:O	1.82	0.97
1:A:87[A]:VAL:HG21	1:A:103:TYR:CE1	2.02	0.93
2:A:606:POY:O16	4:A:1235:HOH:O	1.87	0.93
1:A:151:LYS:HE3	4:A:1133:HOH:O	1.74	0.86
1:A:87[A]:VAL:CG2	1:A:103:TYR:CE1	2.59	0.85
1:A:212:ALA:HB1	2:A:607[A]:POY:H15A	1.57	0.85
1:A:192[A]:PHE:HZ	1:A:570[A]:TRP:CZ3	1.95	0.84
1:A:511[B]:MET:CE	1:A:544:ILE:CD1	2.57	0.82
1:A:87[A]:VAL:CG2	1:A:103:TYR:HE1	1.94	0.81
1:A:244[A]:GLU:OE2	4:A:730:HOH:O	1.99	0.79
1:A:174:ASP:OD1	4:A:1230:HOH:O	1.98	0.79
1:A:277:LEU:CD1	2:A:607[B]:POY:H22	2.12	0.79
1:A:79:ASN:OD1	1:A:89[B]:THR:HG22	1.83	0.78
1:A:277:LEU:HD12	2:A:607[B]:POY:H22	1.65	0.78
1:A:511[B]:MET:CE	1:A:544:ILE:HD11	2.14	0.76
1:A:192[A]:PHE:HZ	1:A:570[A]:TRP:HZ3	1.31	0.76
2:A:607[A]:POY:O18	2:A:607[A]:POY:C11	2.34	0.76
1:A:192[A]:PHE:CZ	1:A:570[A]:TRP:CZ3	2.74	0.75
2:A:607[B]:POY:O17	2:A:607[B]:POY:C20	2.32	0.75

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:ASP:HB2	2:A:607[A]:POY:H13A	1.70	0.72
1:A:192[A]:PHE:HE2	1:A:570[A]:TRP:HH2	1.38	0.72
2:A:606:POY:C14	4:A:1235:HOH:O	2.37	0.71
1:A:192[A]:PHE:CZ	1:A:570[A]:TRP:HZ3	2.08	0.71
1:A:123:LYS:HD3	4:A:1084:HOH:O	1.89	0.70
1:A:511[B]:MET:SD	1:A:544:ILE:HD13	2.32	0.68
2:A:607[A]:POY:C8	2:A:607[A]:POY:H22A	2.22	0.65
1:A:123:LYS:CD	4:A:1084:HOH:O	2.46	0.62
2:A:607[A]:POY:O18	2:A:607[A]:POY:H11	2.00	0.62
1:A:424:ASP:OD1	1:A:428[A]:GLU:HG3	2.01	0.61
1:A:414:ARG:HD2	4:A:895:HOH:O	2.02	0.60
1:A:192[A]:PHE:CE2	1:A:570[A]:TRP:HH2	2.17	0.60
1:A:595:PRO:HA	4:A:1668:HOH:O	2.00	0.60
1:A:370:LYS:HZ3	2:A:606:POY:H15B	1.68	0.59
1:A:431:LYS:HE3	4:A:823:HOH:O	2.02	0.59
1:A:592[B]:VAL:HG12	1:A:593:THR:N	2.16	0.59
1:A:192[A]:PHE:CE2	1:A:570[A]:TRP:CH2	2.89	0.59
1:A:161:MET:CE	1:A:587[A]:MET:HG3	2.33	0.58
1:A:288:LYS:HB3	4:A:767:HOH:O	2.04	0.56
1:A:511[B]:MET:HE2	1:A:544:ILE:HD11	1.87	0.56
1:A:151:LYS:CE	4:A:1133:HOH:O	2.44	0.55
1:A:87[A]:VAL:HG23	1:A:103:TYR:HE1	1.71	0.54
1:A:229:ASP:CA	2:A:607[A]:POY:H15B	2.22	0.53
1:A:208[A]:ASN:OD1	4:A:1126:HOH:O	2.19	0.52
1:A:367:GLY:HA3	2:A:606:POY:H15	1.92	0.52
1:A:89[B]:THR:HG21	4:A:892:HOH:O	2.09	0.52
1:A:140:LYS:HD3	4:A:718:HOH:O	2.09	0.52
1:A:285:VAL:HG13	4:A:1575:HOH:O	2.10	0.50
1:A:288:LYS:HE3	4:A:639:HOH:O	2.10	0.50
1:A:89[A]:THR:HG22	1:A:90:GLY:N	2.26	0.50
1:A:47:THR:HG23	4:A:1104:HOH:O	2.12	0.49
2:A:607[A]:POY:H15A	4:A:1276:HOH:O	1.85	0.48
1:A:511[B]:MET:HE1	1:A:544:ILE:HD11	1.94	0.48
1:A:511[B]:MET:HE1	1:A:544:ILE:CD1	2.42	0.48
1:A:431:LYS:HE2	4:A:661:HOH:O	2.14	0.48
1:A:47:THR:CG2	4:A:1104:HOH:O	2.61	0.47
1:A:288:LYS:HD2	4:A:1088:HOH:O	2.14	0.47
2:A:607[A]:POY:O18	2:A:607[A]:POY:H11A	2.13	0.47
1:A:87[B]:VAL:HG13	4:A:1554:HOH:O	2.14	0.47
1:A:277:LEU:HD11	2:A:607[B]:POY:H22	1.91	0.46
1:A:549:HIS:CE1	1:A:553:LEU:HD11	2.50	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:592[B]:VAL:CG1	1:A:593:THR:N	2.79	0.45
1:A:586:GLY:HA3	4:A:1462:HOH:O	2.16	0.45
1:A:192[A]:PHE:CZ	1:A:570[A]:TRP:CH2	3.04	0.44
1:A:228[A]:HIS:CE1	2:A:607[A]:POY:O18	2.71	0.44
1:A:370:LYS:HZ3	2:A:606:POY:C15	2.30	0.44
1:A:188:ILE:HD12	1:A:266:LEU:HD21	2.00	0.43
1:A:87[A]:VAL:CG2	1:A:103:TYR:CD1	3.00	0.43
3:A:3500:GOL:H11	4:A:696:HOH:O	2.19	0.43
2:A:607[A]:POY:C22	2:A:607[A]:POY:H8	2.22	0.42
1:A:407[A]:LEU:HA	1:A:437:ALA:HB3	2.02	0.42
1:A:161:MET:HE1	1:A:587[A]:MET:HG3	2.03	0.41
2:A:607[A]:POY:C22	2:A:607[A]:POY:C8	2.88	0.41
1:A:305:LYS:HE3	4:A:10:HOH:O	2.20	0.41
1:A:426:ASN:HB2	1:A:428[A]:GLU:HG3	2.03	0.40
1:A:299:LYS:NZ	4:A:733:HOH:O	2.52	0.40
1:A:56:LYS:HD2	4:A:989:HOH:O	2.20	0.40
1:A:79:ASN:OD1	1:A:89[B]:THR:CG2	2.64	0.40
1:A:46:GLU:N	4:A:732:HOH:O	1.63	0.40

All (7) 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)
4:A:1073:HOH:O	4:A:1196:HOH:O[1_565]	1.63	0.57
4:A:1196:HOH:O	4:A:1234:HOH:O[1_545]	1.70	0.50
4:A:1191:HOH:O	4:A:1227:HOH:O[2_545]	1.78	0.42
4:A:963:HOH:O	4:A:1226:HOH:O[2_545]	1.83	0.37
1:A:520:ASN:ND2	4:A:1147:HOH:O[1_565]	1.94	0.26
1:A:244[A]:GLU:OE2	4:A:779:HOH:O[2_556]	2.08	0.12
4:A:1225:HOH:O	4:A:1516:HOH:O[2_556]	2.15	0.05

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	607/584 (104%)	588 (97%)	19 (3%)	0	100 100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	494/469 (105%)	491 (99%)	3 (1%)	90 69

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	LYS
1	A	428[A]	GLU
1	A	428[B]	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	ASN
1	A	447	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

4 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	$\#\|Z\| > 2$
3	GOL	A	3500	-	5,5,5	0.77	0	5,5,5	0.60	0
2	POY	A	606	-	16,23,23	1.82	5 (31%)	16,32,32	2.48	5 (31%)
2	POY	A	607[A]	-	16,23,23	1.56	2 (12%)	16,32,32	1.72	2 (12%)
2	POY	A	607[B]	-	16,23,23	1.50	3 (18%)	16,32,32	3.63	8 (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
3	GOL	A	3500	-	-	0/4/4/4	0/0/0/0
2	POY	A	606	-	-	0/10/10/10	0/2/2/2
2	POY	A	607[A]	-	-	2/10/10/10	0/2/2/2
2	POY	A	607[B]	-	-	0/10/10/10	0/2/2/2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	606	POY	C8-N9	-3.52	1.27	1.34
2	A	606	POY	C15-C14	-2.77	1.39	1.49
2	A	606	POY	C20-N7	-2.65	1.43	1.50
2	A	606	POY	C4-N5	-2.40	1.36	1.39
2	A	607[B]	POY	C2-N1	-2.35	1.34	1.38
2	A	606	POY	O16-C14	2.33	1.29	1.21

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	607[A]	POY	C13-C14	2.82	1.56	1.50
2	A	607[B]	POY	C15-C14	2.94	1.61	1.49
2	A	607[B]	POY	C13-C14	3.80	1.58	1.50
2	A	607[A]	POY	C2-N1	4.55	1.45	1.38

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	607[B]	POY	C3-C2-N1	-7.02	116.38	120.52
2	A	606	POY	C3-C2-N1	-6.52	116.68	120.52
2	A	606	POY	C11-C12-C13	-3.12	101.85	113.29
2	A	606	POY	C11-C10-N1	-3.01	109.75	112.29
2	A	607[B]	POY	C12-C13-C14	-2.47	108.91	116.15
2	A	607[B]	POY	C22-C21-C20	-2.30	92.31	110.66
2	A	607[B]	POY	C12-C11-C10	-2.23	103.90	112.38
2	A	607[B]	POY	C3-C4-N9	-2.07	106.82	110.83
2	A	606	POY	C19-N5-C6	2.94	125.82	119.51
2	A	607[B]	POY	C20-N7-C8	3.20	144.14	126.12
2	A	607[A]	POY	C11-C10-N1	3.83	115.52	112.29
2	A	606	POY	C2-C3-C4	3.96	122.76	119.93
2	A	607[A]	POY	C10-N1-C2	4.39	123.41	119.10
2	A	607[B]	POY	C11-C10-N1	7.46	118.59	112.29
2	A	607[B]	POY	C2-C3-C4	8.20	125.80	119.93

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	607[A]	POY	C11-C10-N1-C6
2	A	607[A]	POY	C11-C10-N1-C2

There are no ring outliers.

4 monomers are involved in 29 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	3500	GOL	1	0
2	A	606	POY	5	0
2	A	607[A]	POY	17	0
2	A	607[B]	POY	6	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	575/584 (98%)	0.14	9 (1%) 74 76	6, 12, 24, 64	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	596	ASN	4.2
1	A	35	ASN	3.6
1	A	70	TRP	3.6
1	A	93	THR	3.0
1	A	71	SER	2.9
1	A	72	GLY	2.3
1	A	34	ASN	2.2
1	A	352	ALA	2.1
1	A	33	SER	2.1

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

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

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

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	POY	A	607[A]	22/22	0.74	0.33	18.80	23,48,56,62	22
2	POY	A	607[B]	22/22	0.74	0.33	15.88	8,16,21,23	22
3	GOL	A	3500	6/6	0.93	0.12	3.62	16,19,20,21	0
2	POY	A	606	22/22	0.89	0.10	2.46	12,19,22,24	0

6.5 Other polymers (i)

There are no such residues in this entry.