



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 23, 2024 – 02:54 PM EDT

PDB ID : 5GZS
Title : Structure of VC protein
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Deposited on : 2016-10-01
Resolution : 2.60 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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	:	2.37.1
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)	:	2.37.1

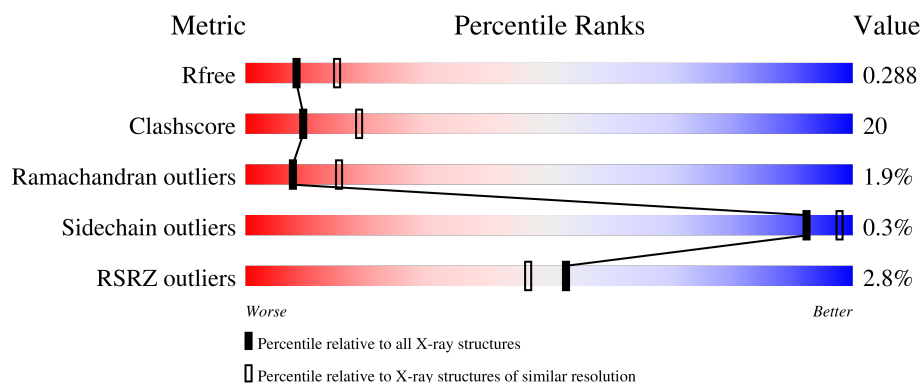
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 $\leq 5\%$. 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	446	

2 Entry composition [i](#)

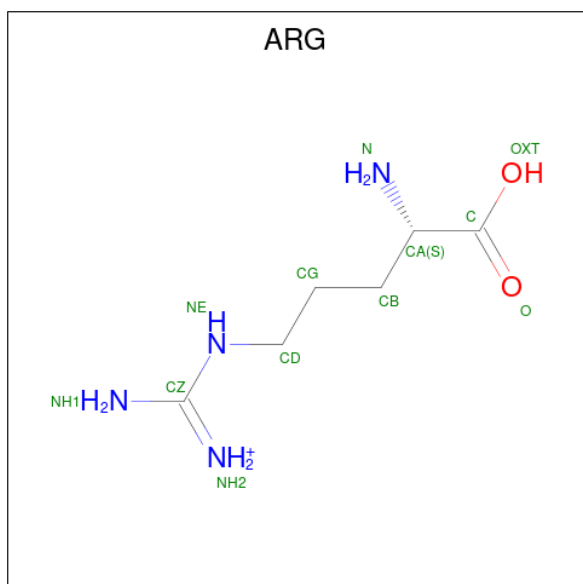
There are 3 unique types of molecules in this entry. The entry contains 3570 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GGDEF family protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	432	Total	C	N	O	S	Se	0	0	0
			3484	2234	583	659	2	6			

- Molecule 2 is ARGinine (three-letter code: ARG) (formula: C₆H₁₅N₄O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			12	6	4	2		

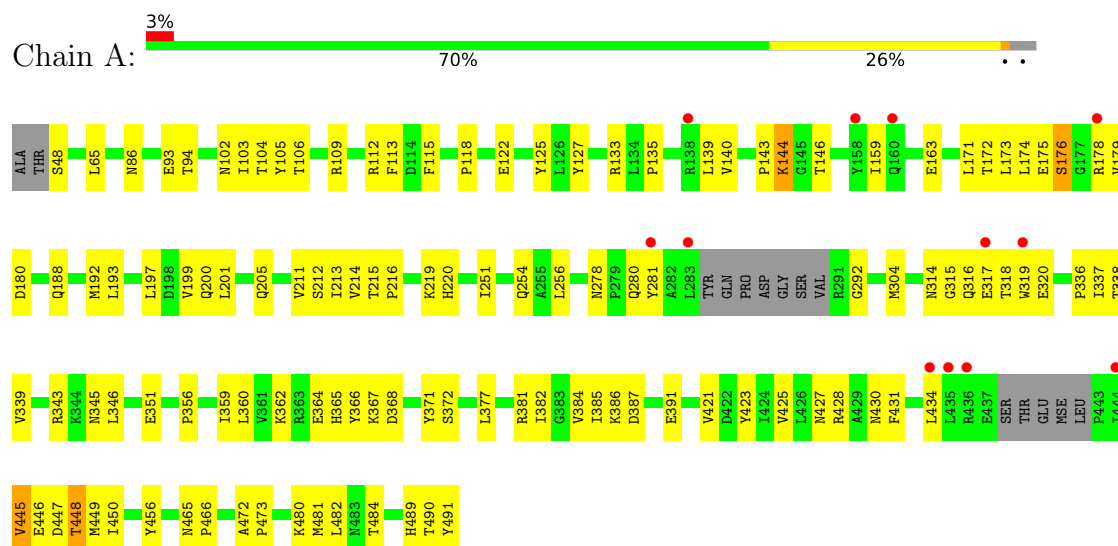
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	74	Total	O	0	0
			74	74		

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: GGDEF family protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	111.71Å 111.71Å 138.65Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.00 – 2.60 39.67 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.8 (40.00-2.60) 95.5 (39.67-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.73 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.219 , 0.287 0.229 , 0.288	Depositor DCC
R_{free} test set	814 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	38.2	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 44.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3570	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.36% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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](#)

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	0.55	0/3544	0.66	1/4794 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	144	LYS	CD-CE-NZ	5.89	125.25	111.70

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	3484	0	3495	137	0
2	A	12	0	12	1	0
3	A	74	0	0	13	0
All	All	3570	0	3507	137	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 20.

All (137) 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:445:VAL:HG22	1:A:446:GLU:H	1.01	1.14
1:A:102:ASN:HA	1:A:211:VAL:CG1	1.90	1.01
1:A:254:GLN:OE1	3:A:601:HOH:O	1.84	0.95
1:A:102:ASN:HA	1:A:211:VAL:HG11	1.44	0.95
1:A:445:VAL:HG22	1:A:446:GLU:N	1.83	0.91
1:A:316:GLN:HB3	1:A:317:GLU:HB3	1.60	0.83
1:A:445:VAL:CG2	1:A:446:GLU:H	1.85	0.83
1:A:173:LEU:HB3	1:A:179:VAL:HG21	1.59	0.82
1:A:113:PHE:CD1	1:A:214:VAL:HG21	2.15	0.82
1:A:104:THR:O	1:A:109:ARG:NH1	2.12	0.81
1:A:427:ASN:OD1	1:A:430:ASN:HB2	1.81	0.81
1:A:316:GLN:HB3	1:A:317:GLU:CA	2.12	0.79
1:A:316:GLN:HB3	1:A:317:GLU:CB	2.13	0.79
1:A:316:GLN:CB	1:A:317:GLU:HB3	2.12	0.79
1:A:316:GLN:HB3	1:A:317:GLU:HA	1.68	0.75
1:A:364:GLU:HG2	1:A:365:HIS:ND1	2.01	0.75
1:A:144:LYS:HB2	1:A:163:GLU:HG3	1.72	0.71
1:A:428:ARG:NH1	1:A:446:GLU:OE2	2.23	0.70
1:A:113:PHE:CE1	1:A:214:VAL:HG21	2.26	0.70
1:A:140:VAL:HG23	1:A:159:ILE:HG21	1.72	0.69
1:A:102:ASN:CA	1:A:211:VAL:CG1	2.69	0.69
1:A:109:ARG:HA	1:A:112:ARG:HD3	1.73	0.69
1:A:484:THR:HG23	3:A:646:HOH:O	1.93	0.68
1:A:133:ARG:HB3	1:A:135:PRO:HD2	1.76	0.67
1:A:343:ARG:HA	1:A:346:LEU:HD12	1.75	0.67
1:A:447:ASP:OD2	1:A:450:ILE:HD12	1.93	0.67
1:A:171:LEU:O	1:A:175:GLU:N	2.25	0.67
1:A:384:VAL:CG1	3:A:604:HOH:O	2.42	0.67
1:A:317:GLU:O	1:A:320:GLU:N	2.29	0.66
1:A:171:LEU:O	1:A:175:GLU:HB2	1.96	0.65
1:A:118:PRO:HA	1:A:212:SER:HB3	1.77	0.65
1:A:384:VAL:HG12	3:A:604:HOH:O	1.96	0.64
1:A:256:LEU:HB2	1:A:481:MSE:HE1	1.80	0.64
1:A:427:ASN:CG	1:A:430:ASN:HB2	2.17	0.64
1:A:371:TYR:O	1:A:449:MSE:HG2	2.00	0.62
1:A:48:SER:CB	3:A:614:HOH:O	2.47	0.62
1:A:113:PHE:CD1	1:A:214:VAL:CG2	2.83	0.61
1:A:102:ASN:CA	1:A:211:VAL:HG11	2.27	0.61
1:A:173:LEU:HB3	1:A:179:VAL:CG2	2.29	0.60
1:A:102:ASN:HA	1:A:211:VAL:HG12	1.79	0.60
1:A:104:THR:HG1	2:A:501:ARG:N	1.98	0.60
1:A:139:LEU:HD13	1:A:178:ARG:HG3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:446:GLU:HG2	1:A:447:ASP:N	2.16	0.59
1:A:304:MSE:HE2	1:A:482:LEU:HD23	1.85	0.58
1:A:281:TYR:CE2	1:A:336:PRO:HG2	2.38	0.58
1:A:122:GLU:HB2	1:A:211:VAL:HG21	1.87	0.57
1:A:106:THR:HG23	1:A:109:ARG:HD2	1.86	0.57
1:A:86:ASN:CB	1:A:144:LYS:HE3	2.35	0.57
1:A:214:VAL:HG22	1:A:215:THR:H	1.70	0.56
1:A:362:LYS:NZ	1:A:368:ASP:OD1	2.35	0.56
1:A:113:PHE:HD1	1:A:214:VAL:CG2	2.18	0.56
1:A:489:HIS:O	1:A:491:TYR:N	2.39	0.56
1:A:316:GLN:HB2	1:A:317:GLU:HB3	1.86	0.56
1:A:109:ARG:O	1:A:112:ARG:HG3	2.05	0.55
1:A:304:MSE:HE2	1:A:482:LEU:CD2	2.36	0.55
1:A:93:GLU:HB3	1:A:113:PHE:HD2	1.71	0.55
1:A:48:SER:N	3:A:614:HOH:O	2.40	0.55
1:A:317:GLU:O	1:A:319:TRP:N	2.41	0.54
1:A:281:TYR:O	1:A:430:ASN:OD1	2.26	0.54
1:A:428:ARG:NH1	1:A:446:GLU:CD	2.61	0.54
1:A:133:ARG:HA	1:A:200:GLN:OE1	2.07	0.54
1:A:214:VAL:HG22	1:A:215:THR:N	2.23	0.54
1:A:171:LEU:O	1:A:175:GLU:CB	2.56	0.53
1:A:386:LYS:HE3	1:A:387:ASP:OD2	2.09	0.52
1:A:102:ASN:C	1:A:211:VAL:CG1	2.78	0.52
1:A:172:THR:O	1:A:175:GLU:O	2.27	0.52
1:A:251:ILE:HG13	3:A:662:HOH:O	2.10	0.52
1:A:427:ASN:ND2	1:A:430:ASN:HB2	2.25	0.52
1:A:216:PRO:O	1:A:219:LYS:HB2	2.09	0.52
1:A:384:VAL:HG13	3:A:604:HOH:O	2.09	0.52
1:A:484:THR:CG2	3:A:646:HOH:O	2.56	0.52
1:A:193:LEU:HD22	1:A:251:ILE:HG22	1.91	0.51
1:A:174:LEU:O	1:A:175:GLU:C	2.49	0.51
1:A:105:TYR:HA	1:A:115:PHE:CE2	2.46	0.51
1:A:427:ASN:ND2	1:A:430:ASN:H	2.08	0.51
1:A:174:LEU:O	1:A:176:SER:N	2.43	0.51
1:A:351:GLU:OE1	1:A:480:LYS:HE2	2.11	0.50
1:A:113:PHE:HD1	1:A:214:VAL:HG21	1.73	0.50
1:A:174:LEU:HD13	1:A:197:LEU:CD2	2.42	0.50
1:A:281:TYR:CZ	1:A:338:THR:HG23	2.46	0.50
1:A:316:GLN:NE2	1:A:320:GLU:O	2.45	0.49
1:A:316:GLN:CB	1:A:317:GLU:CA	2.85	0.49
1:A:384:VAL:HG11	1:A:391:GLU:N	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:THR:HG21	1:A:220:HIS:HB2	1.94	0.49
1:A:356:PRO:HB2	1:A:456:TYR:CE1	2.48	0.48
1:A:65:LEU:HD21	1:A:213:ILE:HD11	1.93	0.48
1:A:281:TYR:CG	1:A:338:THR:CG2	2.96	0.48
1:A:304:MSE:CE	1:A:482:LEU:HD23	2.44	0.48
1:A:109:ARG:O	1:A:112:ARG:CG	2.62	0.48
1:A:278:ASN:O	1:A:280:GLN:HB2	2.14	0.48
1:A:359:ILE:HG21	1:A:428:ARG:NH1	2.29	0.48
1:A:200:GLN:HE21	1:A:205:GLN:NE2	2.12	0.47
1:A:105:TYR:HD2	1:A:115:PHE:CG	2.33	0.47
1:A:188:GLN:O	1:A:192:MSE:HB2	2.14	0.47
1:A:48:SER:HB2	3:A:614:HOH:O	2.12	0.47
1:A:434:LEU:HD21	3:A:644:HOH:O	2.13	0.47
1:A:281:TYR:CE2	1:A:338:THR:HG23	2.50	0.47
1:A:292:GLY:HA2	1:A:491:TYR:CZ	2.50	0.46
1:A:431:PHE:HZ	1:A:445:VAL:HA	1.80	0.46
1:A:430:ASN:O	1:A:434:LEU:HG	2.16	0.46
1:A:173:LEU:HD22	1:A:179:VAL:CG2	2.47	0.45
1:A:139:LEU:O	1:A:180:ASP:HB2	2.17	0.45
1:A:481:MSE:HE2	3:A:602:HOH:O	2.15	0.45
1:A:102:ASN:C	1:A:211:VAL:HG13	2.37	0.45
1:A:384:VAL:HG12	1:A:385:ILE:H	1.81	0.45
1:A:384:VAL:HG12	1:A:385:ILE:N	2.31	0.45
1:A:127:TYR:CD1	1:A:192:MSE:HG2	2.52	0.45
1:A:281:TYR:OH	1:A:337:ILE:HA	2.17	0.45
1:A:366:TYR:CZ	1:A:367:LYS:HD2	2.53	0.44
1:A:109:ARG:HA	1:A:112:ARG:CD	2.44	0.44
1:A:292:GLY:HA2	1:A:491:TYR:CE1	2.52	0.44
1:A:447:ASP:O	1:A:449:MSE:N	2.51	0.44
1:A:193:LEU:HD11	1:A:199:VAL:HG13	2.00	0.44
1:A:317:GLU:C	1:A:320:GLU:H	2.20	0.44
1:A:281:TYR:HE2	1:A:336:PRO:HB2	1.82	0.44
1:A:216:PRO:HB2	1:A:219:LYS:HG3	1.98	0.44
1:A:281:TYR:CE2	1:A:336:PRO:HB2	2.54	0.43
1:A:364:GLU:OE2	1:A:365:HIS:CE1	2.71	0.43
1:A:372:SER:HA	1:A:449:MSE:HG3	2.00	0.43
1:A:356:PRO:HB2	1:A:456:TYR:CZ	2.53	0.43
1:A:281:TYR:CE1	1:A:338:THR:HG23	2.53	0.43
1:A:427:ASN:OD1	1:A:430:ASN:CB	2.60	0.43
1:A:317:GLU:C	1:A:319:TRP:N	2.72	0.43
1:A:472:ALA:HB3	1:A:473:PRO:HD3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:377:LEU:HA	1:A:377:LEU:HD12	1.77	0.42
1:A:465:ASN:HB2	1:A:466:PRO:HD2	2.01	0.42
1:A:447:ASP:O	1:A:448:THR:C	2.57	0.42
1:A:86:ASN:CG	1:A:144:LYS:HE3	2.40	0.42
1:A:345:ASN:OD1	1:A:345:ASN:N	2.51	0.42
1:A:381:ARG:O	1:A:421:VAL:HG22	2.19	0.42
1:A:382:ILE:HA	1:A:423:TYR:O	2.19	0.42
1:A:103:ILE:HG13	1:A:214:VAL:HG11	2.02	0.41
1:A:143:PRO:HB2	1:A:146:THR:CG2	2.50	0.41
1:A:125:TYR:CE2	1:A:201:LEU:HD13	2.55	0.41
1:A:360:LEU:HD21	1:A:425:VAL:HG22	2.03	0.41
1:A:122:GLU:CB	1:A:211:VAL:HG21	2.49	0.41
1:A:481:MSE:CE	3:A:602:HOH:O	2.68	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	426/446 (96%)	388 (91%)	30 (7%)	8 (2%)	8 15

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	318	THR
1	A	445	VAL
1	A	448	THR
1	A	490	THR
1	A	176	SER
1	A	94	THR
1	A	339	VAL
1	A	315	GLY

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	384/389 (99%)	383 (100%)	1 (0%)	92	98

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	314	ASN

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

Mol	Chain	Res	Type
1	A	205	GLN
1	A	254	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 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
2	ARG	A	501	-	10,11,11	0.80	1 (10%)	11,13,13	1.16	2 (18%)

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	ARG	A	501	-	-	1/11/11/11	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	ARG	OXT-C	-2.31	1.23	1.30

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	ARG	OXT-C-O	-2.77	117.81	124.09
2	A	501	ARG	OXT-C-CA	2.16	120.73	113.38

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	ARG	OXT-C-CA-N

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	ARG	1	0

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	426/446 (95%)	-0.07	12 (2%) 53 46	26, 55, 84, 165	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	281	TYR	6.8
1	A	178	ARG	5.6
1	A	435	LEU	4.2
1	A	158	TYR	3.3
1	A	319	TRP	3.3
1	A	160	GLN	2.8
1	A	436	ARG	2.6
1	A	283	LEU	2.5
1	A	434	LEU	2.4
1	A	138	ARG	2.3
1	A	444	ILE	2.3
1	A	317	GLU	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 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(\AA^2)	Q<0.9
2	ARG	A	501	12/12	0.96	0.20	40,53,70,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.