Full wwPDB X-ray Structure Validation Report (i)

Nov 6, 2023 – 05:12 PM EST

PDB ID	:	6URU
Title	:	iAChSnFR Fluorescent Acetylcholine Sensor precursor
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Deposited on	:	2019-10-24
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 (i) 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 (1)) were used in the production of this report:

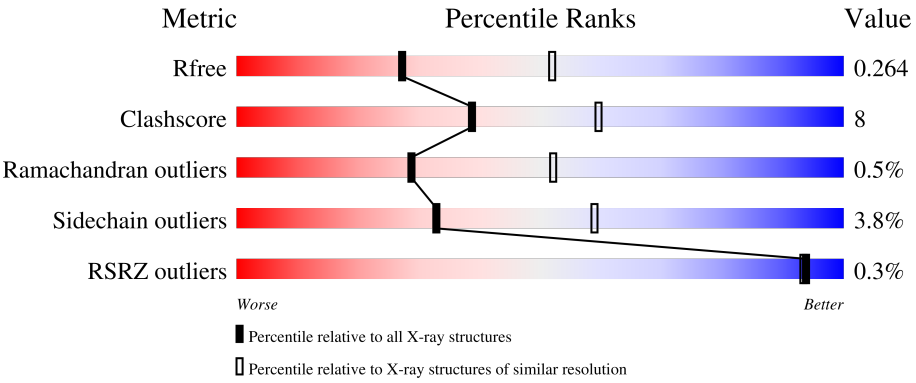
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.60 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 <=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	558	76% 13% • 10%
1	B	558	76% 13% • 10%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 7974 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 iAChSnFR precursor.

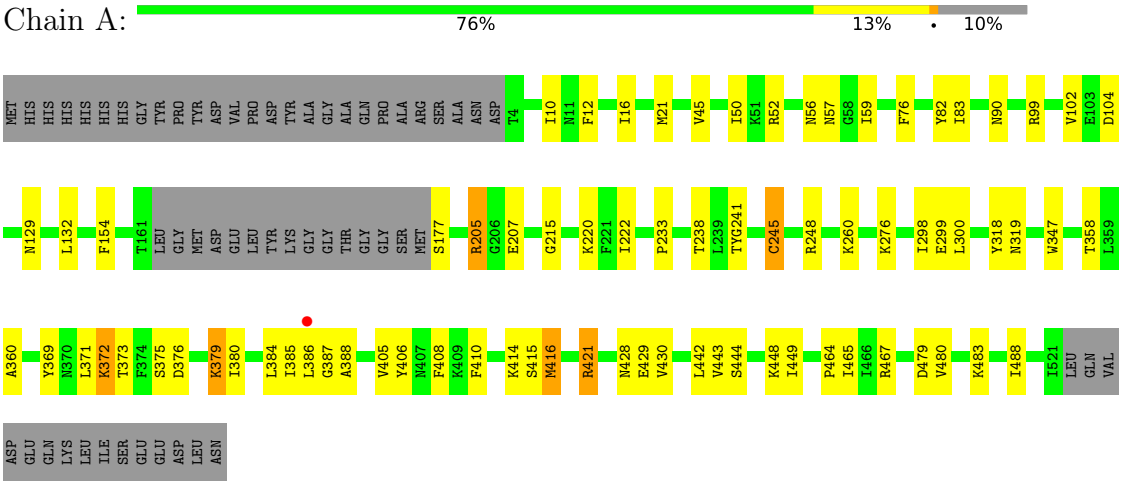
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	501	Total	C	N	O	S	0	0	0
			3987	2554	660	762	11			
1	B	501	Total	C	N	O	S	0	0	0
			3987	2554	660	762	11			

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.

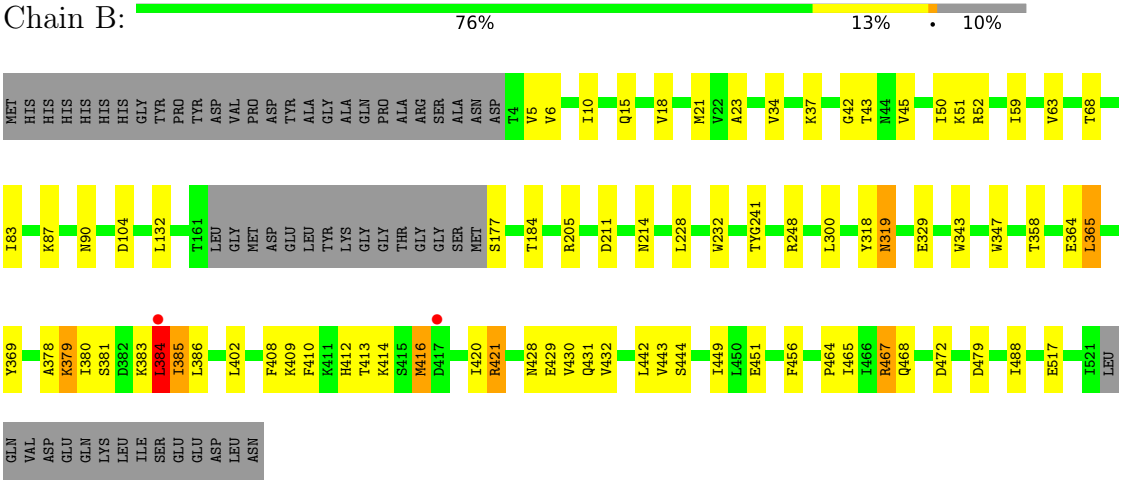
• Molecule 1: iAChSnFR precursor

Chain A:



• Molecule 1: iAChSnFR precursor

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	72.12Å 89.78Å 92.89Å 90.00° 90.48° 90.00°	Depositor
Resolution (Å)	92.89 – 2.60 92.89 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.1 (92.89-2.60) 97.0 (92.89-2.60)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.82 (at 2.62Å)	Xtriage
Refinement program	REFMAC 5.8.0222	Depositor
R, R_{free}	0.213 , 0.265 0.216 , 0.264	Depositor DCC
R_{free} test set	1724 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å^2)	50.9	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 18.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.016 for -h,l,k 0.019 for -h,-l,-k 0.457 for h,-k,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	7974	wwPDB-VP
Average B, all atoms (Å^2)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.18% 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: CRO

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	0.59	0/4049	0.74	0/5477
1	B	0.58	0/4049	0.74	0/5477
All	All	0.58	0/8098	0.74	0/10954

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	0	3
1	B	0	5
All	All	0	8

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	205	ARG	Sidechain
1	A	248	ARG	Sidechain
1	A	52	ARG	Sidechain
1	B	205	ARG	Sidechain
1	B	248	ARG	Sidechain
1	B	421	ARG	Sidechain
1	B	467	ARG	Sidechain
1	B	52	ARG	Sidechain

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	3987	0	3967	51	0
1	B	3987	0	3968	69	0
All	All	7974	0	7935	120	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 8.

All (120) 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:B:385:ILE:HD13	1:B:413:THR:HA	1.41	1.02
1:A:386:LEU:HB3	1:A:430:VAL:HB	1.48	0.95
1:B:383:LYS:O	1:B:384:LEU:O	1.95	0.82
1:A:373:THR:HG22	1:A:376:ASP:OD2	1.79	0.82
1:A:347:TRP:CD2	1:A:465:ILE:HD11	2.16	0.81
1:B:384:LEU:HD23	1:B:431:GLN:OE1	1.80	0.81
1:B:384:LEU:HG	1:B:385:ILE:N	1.96	0.80
1:B:442:LEU:HD13	1:B:449:ILE:HD11	1.64	0.80
1:B:21:MET:CE	1:B:488:ILE:HD11	2.12	0.79
1:B:385:ILE:HD13	1:B:413:THR:CA	2.11	0.79
1:A:386:LEU:HD21	1:A:416:MET:SD	2.23	0.79
1:B:385:ILE:CD1	1:B:413:THR:HA	2.13	0.79
1:B:384:LEU:HD21	1:B:429:GLU:O	1.83	0.79
1:B:21:MET:HE1	1:B:488:ILE:HD11	1.66	0.78
1:B:385:ILE:HG23	1:B:430:VAL:HG12	1.67	0.77
1:B:347:TRP:CD2	1:B:465:ILE:HD11	2.22	0.74
1:B:385:ILE:HD12	1:B:386:LEU:N	2.04	0.72
1:B:386:LEU:O	1:B:432:VAL:O	2.06	0.72
1:A:347:TRP:CE3	1:A:465:ILE:HD11	2.26	0.71
1:A:358:THR:HG21	1:A:449:ILE:HG13	1.73	0.70
1:B:347:TRP:CE3	1:B:465:ILE:HD11	2.26	0.70
1:A:369:TYR:HB2	1:A:371:LEU:HD21	1.74	0.69
1:A:21:MET:HE1	1:A:488:ILE:HD11	1.73	0.69
1:B:442:LEU:HD13	1:B:449:ILE:CD1	2.23	0.67
1:B:365:LEU:HD22	1:B:365:LEU:C	2.17	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:VAL:CG1	1:B:63:VAL:HG21	2.28	0.64
1:B:300:LEU:C	1:B:300:LEU:HD23	2.18	0.63
1:B:408:PHE:CD1	1:B:410:PHE:CE2	2.86	0.63
1:A:442:LEU:HD13	1:A:449:ILE:CD1	2.29	0.63
1:A:386:LEU:O	1:A:416:MET:HG3	1.99	0.63
1:B:10:ILE:HD11	1:B:45:VAL:HG13	1.82	0.62
1:B:318:TYR:O	1:B:318:TYR:CD2	2.54	0.61
1:B:384:LEU:CD2	1:B:429:GLU:O	2.48	0.60
1:B:384:LEU:CD1	1:B:385:ILE:HG22	2.32	0.59
1:A:442:LEU:HD13	1:A:449:ILE:HD11	1.84	0.59
1:A:300:LEU:C	1:A:300:LEU:HD23	2.23	0.59
1:B:442:LEU:CD1	1:B:449:ILE:HD11	2.32	0.59
1:B:385:ILE:HD12	1:B:386:LEU:H	1.68	0.58
1:B:384:LEU:HD21	1:B:430:VAL:HA	1.85	0.58
1:B:381:SER:HA	1:B:409:LYS:O	2.03	0.58
1:B:416:MET:HG3	1:B:420:ILE:HD11	1.85	0.57
1:B:378:ALA:O	1:B:380:ILE:N	2.37	0.57
1:B:21:MET:HE2	1:B:488:ILE:HD11	1.86	0.57
1:A:386:LEU:CD2	1:A:416:MET:SD	2.93	0.56
1:B:364:GLU:HG3	1:B:365:LEU:N	2.21	0.56
1:A:102:VAL:O	1:A:104:ASP:O	2.24	0.55
1:B:386:LEU:H	1:B:386:LEU:HD12	1.73	0.54
1:A:387:GLY:O	1:A:388:ALA:C	2.45	0.54
1:A:408:PHE:CD1	1:A:410:PHE:CE1	2.95	0.54
1:A:371:LEU:O	1:A:373:THR:N	2.38	0.54
1:B:18:VAL:HG12	1:B:63:VAL:HG21	1.90	0.53
1:A:386:LEU:HB3	1:A:430:VAL:CB	2.32	0.53
1:A:464:PRO:O	1:A:465:ILE:HD12	2.09	0.53
1:B:383:LYS:O	1:B:384:LEU:C	2.48	0.52
1:A:50:ILE:O	1:A:467:ARG:NH1	2.41	0.52
1:A:414:LYS:HA	1:A:414:LYS:HE2	1.92	0.51
1:A:318:TYR:CD2	1:A:318:TYR:O	2.64	0.51
1:A:408:PHE:CD1	1:A:410:PHE:HE1	2.29	0.51
1:B:408:PHE:CD1	1:B:410:PHE:HE2	2.29	0.50
1:A:373:THR:HG22	1:A:376:ASP:CG	2.32	0.50
1:B:23:ALA:HB1	1:B:34:VAL:HG11	1.94	0.50
1:B:384:LEU:CD2	1:B:431:GLN:OE1	2.57	0.50
1:B:10:ILE:HG22	1:B:15:GLN:HG3	1.94	0.50
1:B:464:PRO:O	1:B:465:ILE:HD12	2.12	0.50
1:B:358:THR:HG21	1:B:449:ILE:HG13	1.94	0.50
1:A:21:MET:HE1	1:A:488:ILE:CD1	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:ASP:OD2	1:B:214:ASN:ND2	2.46	0.49
1:A:10:ILE:HD11	1:A:45:VAL:CG2	2.43	0.49
1:B:402:LEU:HD13	1:B:456:PHE:CD2	2.48	0.48
1:B:364:GLU:CG	1:B:365:LEU:N	2.76	0.48
1:A:386:LEU:O	1:A:414:LYS:O	2.31	0.48
1:A:369:TYR:HB2	1:A:371:LEU:CD2	2.43	0.48
1:B:365:LEU:HD22	1:B:365:LEU:O	2.14	0.48
1:A:464:PRO:C	1:A:465:ILE:HD12	2.33	0.47
1:A:360:ALA:HA	1:A:448:LYS:O	2.14	0.47
1:A:442:LEU:CD1	1:A:449:ILE:HD11	2.45	0.47
1:A:443:VAL:O	1:A:444:SER:C	2.52	0.47
1:B:43:THR:HG21	1:B:68:THR:HG21	1.97	0.47
1:B:378:ALA:C	1:B:380:ILE:H	2.18	0.47
1:A:154:PHE:HA	1:A:215:GLY:O	2.15	0.46
1:A:238:THR:CG2	1:A:298:ILE:HG21	2.46	0.46
1:A:385:ILE:HD11	1:A:410:PHE:CD1	2.50	0.46
1:A:50:ILE:HA	1:A:59:ILE:O	2.16	0.46
1:B:384:LEU:CG	1:B:385:ILE:N	2.71	0.46
1:A:405:VAL:HG23	1:A:406:TYR:CD2	2.50	0.46
1:A:12:PHE:O	1:A:16:ILE:HG12	2.16	0.45
1:B:42:GLY:O	1:B:45:VAL:HG12	2.16	0.45
1:A:379:LYS:O	1:A:380:ILE:HG13	2.17	0.44
1:A:442:LEU:HD13	1:A:449:ILE:HD13	1.99	0.44
1:A:83:ILE:HD13	1:A:132:LEU:HG	1.99	0.44
1:B:83:ILE:HD13	1:B:132:LEU:HG	1.98	0.44
1:A:76:PHE:CD1	1:A:99:ARG:HD3	2.52	0.44
1:A:480:VAL:O	1:A:483:LYS:HB2	2.18	0.43
1:B:37:LYS:HB3	1:B:59:ILE:HD11	2.00	0.43
1:B:416:MET:HG2	1:B:421:ARG:HA	2.00	0.43
1:B:468:GLN:NE2	1:B:472:ASP:OD1	2.47	0.43
1:B:379:LYS:O	1:B:380:ILE:HG13	2.19	0.43
1:A:387:GLY:O	1:A:421:ARG:HD3	2.18	0.43
1:A:384:LEU:HD13	1:A:429:GLU:O	2.18	0.43
1:A:479:ASP:O	1:A:483:LYS:HG3	2.19	0.43
1:B:464:PRO:C	1:B:465:ILE:HD12	2.38	0.43
1:B:21:MET:HE1	1:B:488:ILE:CD1	2.44	0.43
1:B:10:ILE:HD11	1:B:45:VAL:CG1	2.48	0.42
1:B:443:VAL:O	1:B:444:SER:C	2.57	0.42
1:B:50:ILE:O	1:B:467:ARG:NH1	2.49	0.42
1:B:386:LEU:HD11	1:B:410:PHE:CD2	2.53	0.42
1:B:51:LYS:HE2	1:B:343:TRP:CZ2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:410:PHE:HE1	1:B:432:VAL:HG21	1.85	0.42
1:A:385:ILE:HD11	1:A:410:PHE:CE1	2.55	0.42
1:B:50:ILE:HA	1:B:59:ILE:O	2.20	0.42
1:B:385:ILE:HG12	1:B:414:LYS:CB	2.49	0.42
1:B:43:THR:CG2	1:B:68:THR:HG21	2.49	0.42
1:A:245:CYS:O	1:A:260:LYS:NZ	2.46	0.41
1:A:205:ARG:NH2	1:A:207:GLU:OE2	2.49	0.41
1:B:228:LEU:HD22	1:B:232:TRP:CE2	2.56	0.41
1:A:82:TYR:CZ	1:A:129:ASN:HB3	2.56	0.40
1:B:365:LEU:C	1:B:365:LEU:CD2	2.87	0.40
1:A:220:LYS:HE2	1:A:222:ILE:HD11	2.03	0.40
1:B:6:VAL:HG12	1:B:59:ILE:HD13	2.03	0.40
1:B:83:ILE:HD11	1:B:132:LEU:HD12	2.02	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	494/558 (88%)	465 (94%)	27 (6%)	2 (0%)	34	57
1	B	494/558 (88%)	462 (94%)	29 (6%)	3 (1%)	25	47
All	All	988/1116 (88%)	927 (94%)	56 (6%)	5 (0%)	29	52

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	384	LEU
1	A	379	LYS
1	B	379	LYS
1	B	319	ASN
1	A	372	LYS

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	436/482 (90%)	421 (97%)	15 (3%)	37 63
1	B	436/482 (90%)	418 (96%)	18 (4%)	30 56
All	All	872/964 (90%)	839 (96%)	33 (4%)	33 59

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	ASN
1	A	57	ASN
1	A	90	ASN
1	A	177	SER
1	A	233	PRO
1	A	245	CYS
1	A	276	LYS
1	A	299	GLU
1	A	319	ASN
1	A	372	LYS
1	A	375	SER
1	A	415	SER
1	A	416	MET
1	A	421	ARG
1	A	428	ASN
1	B	5	VAL
1	B	87	LYS
1	B	90	ASN
1	B	104	ASP
1	B	177	SER
1	B	184	THR
1	B	319	ASN
1	B	329	GLU
1	B	365	LEU
1	B	369	TYR
1	B	384	LEU
1	B	385	ILE

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Mol	Chain	Res	Type
1	B	412	HIS
1	B	416	MET
1	B	428	ASN
1	B	451	GLU
1	B	479	ASP
1	B	517	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	344	ASN
1	B	56	ASN
1	B	80	ASN
1	B	214	ASN
1	B	368	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)

2 non-standard protein/DNA/RNA residues 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	$\#\|Z\| > 2$
1	CRO	A	241	1	23,23,24	3.39	6 (26%)	30,32,34	3.95	9 (30%)
1	CRO	B	241	1	23,23,24	3.19	7 (30%)	30,32,34	3.33	15 (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
1	CRO	A	241	1	-	1/12/31/32	0/2/2/2
1	CRO	B	241	1	-	1/12/31/32	0/2/2/2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	241	CRO	CB2-CA2	13.69	1.46	1.35
1	B	241	CRO	CB2-CA2	13.00	1.46	1.35
1	A	241	CRO	CA2-C2	-5.40	1.43	1.48
1	B	241	CRO	O2-C2	4.27	1.32	1.23
1	A	241	CRO	O2-C2	3.52	1.30	1.23
1	B	241	CRO	C1-N2	3.41	1.37	1.32
1	A	241	CRO	C1-N2	3.33	1.37	1.32
1	B	241	CRO	CA2-C2	-3.06	1.45	1.48
1	A	241	CRO	CA2-N2	-2.84	1.32	1.38
1	A	241	CRO	C2-N3	-2.75	1.33	1.39
1	B	241	CRO	CA1-C1	-2.70	1.47	1.51
1	B	241	CRO	CA2-N2	-2.41	1.33	1.38
1	B	241	CRO	C2-N3	-2.06	1.35	1.39

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	241	CRO	O2-C2-CA2	-16.92	121.46	130.96
1	B	241	CRO	O2-C2-CA2	-13.00	123.66	130.96
1	A	241	CRO	CA2-C2-N3	8.76	107.51	103.37
1	B	241	CRO	O3-C3-CA3	-6.49	106.79	126.39
1	A	241	CRO	O3-C3-CA3	-6.26	107.49	126.39
1	B	241	CRO	CA2-C2-N3	5.70	106.06	103.37
1	A	241	CRO	O2-C2-N3	3.35	131.00	124.35
1	B	241	CRO	O2-C2-N3	2.98	130.28	124.35
1	B	241	CRO	CG1-CB1-CA1	-2.89	105.34	112.16
1	B	241	CRO	C2-N3-C1	-2.65	106.62	107.97
1	B	241	CRO	CA1-C1-N3	-2.51	121.74	124.75
1	B	241	CRO	CA3-N3-C1	-2.49	124.18	127.16
1	B	241	CRO	CE1-CD1-CG2	-2.47	118.03	121.25
1	B	241	CRO	CD2-CG2-CB2	-2.38	113.11	121.22
1	A	241	CRO	CD2-CG2-CD1	2.35	121.11	117.64
1	A	241	CRO	CG2-CB2-CA2	-2.34	127.07	129.94
1	A	241	CRO	CA1-C1-N2	2.33	127.15	123.89
1	B	241	CRO	CD2-CG2-CD1	2.31	121.06	117.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	241	CRO	CA3-N3-C2	2.25	128.96	123.80
1	B	241	CRO	CB2-CA2-N2	-2.22	125.74	128.83
1	B	241	CRO	CG2-CB2-CA2	-2.06	127.42	129.94
1	A	241	CRO	CD2-CG2-CB2	-2.02	114.35	121.22
1	B	241	CRO	N3-C1-N2	2.01	112.84	111.45
1	A	241	CRO	CE1-CD1-CG2	-2.01	118.63	121.25

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	241	CRO	C3-CA3-N3-C2
1	A	241	CRO	C3-CA3-N3-C2

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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

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	500/558 (89%)	-0.37	1 (0%)	95	95	32, 51, 90, 141	0
1	B	500/558 (89%)	-0.36	2 (0%)	92	91	31, 51, 86, 152	0
All	All	1000/1116 (89%)	-0.36	3 (0%)	94	93	31, 51, 88, 152	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	386	LEU	6.2
1	B	384	LEU	3.0
1	B	417	ASP	3.0

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

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, 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	B-factors(Å^2)	Q<0.9
1	CRO	B	241	22/23	0.95	0.16	36,42,47,51	0
1	CRO	A	241	22/23	0.96	0.17	35,45,51,52	0

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.