Full wwPDB X-ray Structure Validation Report (i)

Jun 25, 2024 – 09:01 AM EDT

PDB ID	:	5TXE
Title	:	AtxE2 Isopeptidase - S527A Variant with Astexin3-dC4 Bound
Authors	:	Chekan, J.R.; Nair, S.K.
Deposited on	:	2016-11-16
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 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:

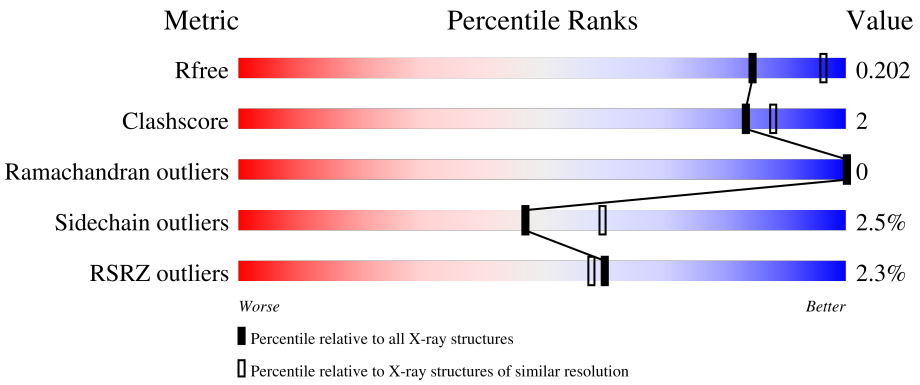
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (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	705	3%, 86%, 5% • 8%
1	B	705	%, 84%, 8% • 8%
2	C	24	46%, 12%, 8% • 29%
2	D	24	63%, 8% • 25%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 11501 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 AtxE2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	652	5161	3248	935	963	15	0	3	0
1	B	652	5171	3255	937	964	15	0	4	0

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	527	ALA	SER	engineered mutation	UNP E8RUP5
A	696	GLY	-	expression tag	UNP E8RUP5
A	697	SER	-	expression tag	UNP E8RUP5
A	698	ARG	-	expression tag	UNP E8RUP5
A	699	SER	-	expression tag	UNP E8RUP5
A	700	HIS	-	expression tag	UNP E8RUP5
A	701	HIS	-	expression tag	UNP E8RUP5
A	702	HIS	-	expression tag	UNP E8RUP5
A	703	HIS	-	expression tag	UNP E8RUP5
A	704	HIS	-	expression tag	UNP E8RUP5
A	705	HIS	-	expression tag	UNP E8RUP5
B	527	ALA	SER	engineered mutation	UNP E8RUP5
B	696	GLY	-	expression tag	UNP E8RUP5
B	697	SER	-	expression tag	UNP E8RUP5
B	698	ARG	-	expression tag	UNP E8RUP5
B	699	SER	-	expression tag	UNP E8RUP5
B	700	HIS	-	expression tag	UNP E8RUP5
B	701	HIS	-	expression tag	UNP E8RUP5
B	702	HIS	-	expression tag	UNP E8RUP5
B	703	HIS	-	expression tag	UNP E8RUP5
B	704	HIS	-	expression tag	UNP E8RUP5
B	705	HIS	-	expression tag	UNP E8RUP5

- Molecule 2 is a protein called Astexin3-dC4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	17	125	80	19	25	1	0	0	0
			Total	C	N	O	S			
2	D	18	134	85	21	27	1	0	0	0

- Molecule 3 is water.

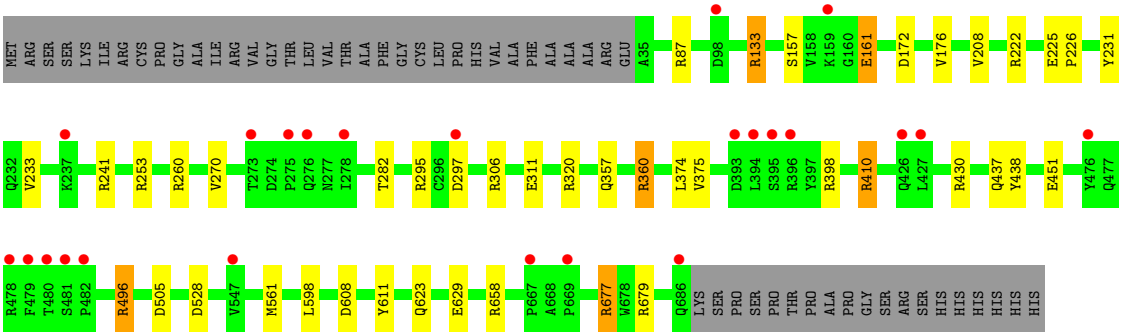
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	339	339	339	0	0
			Total	O		
3	C	10	10	10	0	0
			Total	O		
3	B	551	551	551	0	0
			Total	O		
3	D	10	10	10	0	0

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.

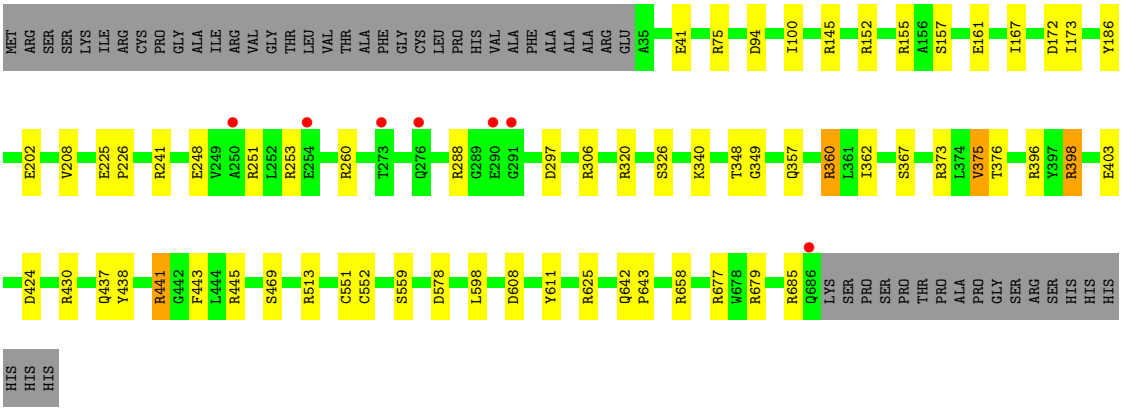
- Molecule 1: AtxE2

Chain A:



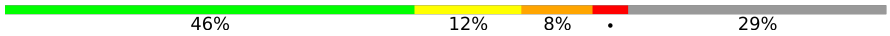
- Molecule 1: AtxE2

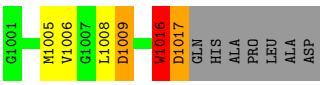
Chain B:



- Molecule 2: Astexin3-dC4

Chain C:





- Molecule 2: Astexin3-dC4

Chain D:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	151.25Å 202.85Å 110.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.82 – 2.20 29.82 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.82-2.20) 99.9 (29.82-2.20)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.87 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.154 , 0.198 0.163 , 0.202	Depositor DCC
R_{free} test set	4255 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å^2)	29.8	Xtriage
Anisotropy	0.272	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 39.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	11501	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% 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)

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	0.82	1/5297 (0.0%)	0.93	13/7224 (0.2%)
1	B	0.92	2/5304 (0.0%)	1.06	32/7233 (0.4%)
2	C	0.91	0/129	1.21	2/176 (1.1%)
2	D	0.90	0/138	0.83	0/188
All	All	0.87	3/10868 (0.0%)	1.00	47/14821 (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
2	C	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	241	ARG	CD-NE	-6.48	1.35	1.46
1	A	451	GLU	CD-OE1	5.27	1.31	1.25
1	B	559	SER	CB-OG	-5.05	1.35	1.42

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	445	ARG	NE-CZ-NH2	-21.03	109.78	120.30
1	B	445	ARG	NE-CZ-NH1	15.88	128.24	120.30
1	B	241	ARG	NE-CZ-NH2	-12.03	114.28	120.30
1	A	677	ARG	NE-CZ-NH2	-10.42	115.09	120.30
1	A	677	ARG	NE-CZ-NH1	9.65	125.13	120.30
1	B	679	ARG	NE-CZ-NH2	-9.63	115.48	120.30
1	B	679	ARG	NE-CZ-NH1	9.19	124.89	120.30
1	B	441	ARG	NE-CZ-NH2	-8.71	115.95	120.30
1	A	222	ARG	NE-CZ-NH1	8.62	124.61	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	445	ARG	CD-NE-CZ	8.57	135.60	123.60
1	B	241	ARG	NE-CZ-NH1	8.32	124.46	120.30
1	B	260	ARG	NE-CZ-NH1	7.66	124.13	120.30
1	B	253	ARG	NE-CZ-NH1	7.17	123.89	120.30
1	B	430	ARG	NE-CZ-NH1	6.80	123.70	120.30
1	B	360	ARG	NE-CZ-NH1	6.75	123.68	120.30
1	B	360	ARG	NE-CZ-NH2	-6.73	116.94	120.30
2	C	1016	TRP	CB-CA-C	6.70	123.79	110.40
1	B	373	ARG	NE-CZ-NH2	-6.69	116.95	120.30
1	B	441	ARG	CG-CD-NE	-6.52	98.11	111.80
1	A	679	ARG	NE-CZ-NH2	-6.40	117.10	120.30
1	B	152	ARG	NE-CZ-NH1	6.36	123.48	120.30
1	B	625	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	A	679	ARG	NE-CZ-NH1	6.23	123.42	120.30
1	B	685	ARG	NE-CZ-NH1	6.22	123.41	120.30
1	A	133	ARG	CG-CD-NE	6.20	124.81	111.80
2	C	1016	TRP	CA-C-N	6.11	130.64	117.20
1	A	320	ARG	NE-CZ-NH2	-5.92	117.34	120.30
1	B	320	ARG	NE-CZ-NH2	-5.82	117.39	120.30
1	A	430	ARG	NE-CZ-NH1	5.70	123.15	120.30
1	B	513	ARG	NE-CZ-NH1	5.70	123.15	120.30
1	B	152	ARG	NE-CZ-NH2	-5.61	117.49	120.30
1	B	375	VAL	CG1-CB-CG2	5.57	119.81	110.90
1	A	608	ASP	CB-CG-OD1	5.50	123.25	118.30
1	B	260	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	B	513	ARG	NE-CZ-NH2	-5.41	117.60	120.30
1	A	222	ARG	NE-CZ-NH2	-5.32	117.64	120.30
1	B	145	ARG	NE-CZ-NH2	-5.27	117.66	120.30
1	B	251	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	B	398	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	B	608	ASP	CB-CG-OD2	-5.18	113.64	118.30
1	A	172	ASP	CB-CG-OD2	-5.17	113.65	118.30
1	A	658	ARG	NE-CZ-NH2	-5.14	117.73	120.30
1	B	608	ASP	CB-CG-OD1	5.10	122.89	118.30
1	B	94	ASP	CB-CG-OD1	5.07	122.86	118.30
1	B	172	ASP	CB-CG-OD1	5.03	122.83	118.30
1	B	658	ARG	NE-CZ-NH1	5.02	122.81	120.30
1	A	561	MET	CG-SD-CE	-5.02	92.17	100.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	1016	TRP	Peptide

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	5161	0	5068	18	0
1	B	5171	0	5080	26	1
2	C	125	0	112	6	0
2	D	134	0	120	4	0
3	A	339	0	0	2	0
3	B	551	0	0	11	0
3	C	10	0	0	0	0
3	D	10	0	0	0	0
All	All	11501	0	10380	51	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 2.

All (51) 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)
2:C:1016:TRP:CD1	2:C:1017:ASP:HA	1.68	1.27
1:B:348:THR:C	3:B:802:HOH:O	1.91	1.07
2:D:1016:TRP:CD1	2:D:1017:ASP:N	2.22	1.07
2:C:1016:TRP:CD1	2:C:1017:ASP:CA	2.55	0.88
1:B:348:THR:CB	3:B:802:HOH:O	2.26	0.83
1:B:202[B]:GLU:OE2	3:B:801:HOH:O	2.00	0.79
1:A:410:ARG:NH1	1:A:505:ASP:OD1	2.20	0.75
1:B:41:GLU:HG3	3:B:1067:HOH:O	1.89	0.71
1:A:157:SER:OG	3:A:801:HOH:O	2.09	0.69
2:D:1016:TRP:CG	2:D:1017:ASP:N	2.61	0.64
1:A:270:VAL:HG13	1:A:282:THR:HB	1.80	0.64
1:B:326:SER:O	1:B:441:ARG:NH2	2.31	0.59
1:B:398:ARG:NH1	1:B:424:ASP:OD2	2.36	0.58
1:B:362:ILE:HD13	1:B:376:THR:HG22	1.87	0.57
1:B:155:ARG:NH2	3:B:810:HOH:O	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:GLU:CG	3:B:1067:HOH:O	2.50	0.56
2:C:1016:TRP:NE1	2:C:1017:ASP:HA	2.16	0.55
1:A:133:ARG:NH1	1:A:176:VAL:O	2.40	0.55
1:A:437:GLN:HB3	1:A:438:TYR:HB3	1.88	0.55
1:A:629:GLU:OE2	1:A:677:ARG:HD2	2.07	0.54
2:D:1016:TRP:HD1	2:D:1017:ASP:N	2.01	0.54
2:D:1009:ASP:OD1	2:D:1016:TRP:HB2	2.08	0.53
1:B:349:GLY:N	3:B:802:HOH:O	2.25	0.52
1:B:437:GLN:HB3	1:B:438:TYR:HB3	1.92	0.51
1:B:357:GLN:OE1	1:B:360:ARG:NH1	2.40	0.51
1:A:496:ARG:HH22	1:A:528:ASP:CG	2.16	0.49
2:C:1009:ASP:OD1	2:C:1016:TRP:HB2	2.12	0.49
1:A:157:SER:OG	1:A:161:GLU:HG2	2.14	0.47
1:A:225:GLU:OE1	1:A:226:PRO:HA	2.15	0.47
1:B:225:GLU:OE1	1:B:226:PRO:HA	2.15	0.47
1:A:231:TYR:OH	1:A:253:ARG:NH1	2.47	0.46
1:A:677:ARG:CZ	1:B:208:VAL:HG21	2.46	0.46
1:B:173[A]:ILE:HD12	1:B:186:TYR:CD2	2.52	0.44
1:A:233:VAL:O	1:A:241:ARG:HA	2.17	0.44
1:A:157:SER:CB	1:A:161:GLU:HG2	2.48	0.44
2:C:1006:VAL:HG12	2:C:1017:ASP:HB2	2.01	0.43
1:B:75:ARG:NH2	3:B:831:HOH:O	2.52	0.43
1:B:167:ILE:HG22	1:B:173[B]:ILE:HD11	2.01	0.43
1:B:100:ILE:HD11	3:B:1206:HOH:O	2.19	0.42
1:A:357:GLN:OE1	1:A:360:ARG:NH1	2.48	0.42
1:A:208:VAL:HG21	1:B:677:ARG:CZ	2.50	0.42
1:A:133:ARG:HA	1:A:133:ARG:HD3	1.84	0.42
1:B:248:GLU:HA	3:B:873:HOH:O	2.19	0.41
1:A:87:ARG:NH2	3:A:829:HOH:O	2.53	0.41
1:A:623:GLN:HE22	1:B:578:ASP:HA	1.84	0.41
1:B:642:GLN:HA	1:B:643:PRO:HD3	1.95	0.41
1:B:367:SER:O	1:B:443:PHE:HA	2.21	0.41
2:C:1016:TRP:HD1	2:C:1017:ASP:CA	2.24	0.40
1:B:438:TYR:HB2	1:B:469:SER:HB2	2.03	0.40
1:B:288:ARG:HD3	3:B:1163:HOH:O	2.22	0.40
1:B:551:CYS:HA	1:B:552:CYS:HA	1.88	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:B:403:GLU:OE2	1:B:403:GLU:OE2[6_565]	2.18	0.02

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	653/705 (93%)	640 (98%)	13 (2%)	0	100	100
1	B	654/705 (93%)	640 (98%)	14 (2%)	0	100	100
2	C	15/24 (62%)	12 (80%)	3 (20%)	0	100	100
2	D	16/24 (67%)	12 (75%)	4 (25%)	0	100	100
All	All	1338/1458 (92%)	1304 (98%)	34 (2%)	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	548/587 (93%)	533 (97%)	15 (3%)	44	57
1	B	549/587 (94%)	540 (98%)	9 (2%)	62	76
2	C	14/19 (74%)	10 (71%)	4 (29%)	0	0
2	D	15/19 (79%)	14 (93%)	1 (7%)	16	18
All	All	1126/1212 (93%)	1097 (97%)	29 (3%)	47	58

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	161	GLU
1	A	260	ARG
1	A	295	ARG
1	A	297[A]	ASP
1	A	297[B]	ASP
1	A	306	ARG
1	A	311	GLU
1	A	360	ARG
1	A	374	LEU
1	A	375	VAL
1	A	398	ARG
1	A	410	ARG
1	A	496	ARG
1	A	598	LEU
1	A	611	TYR
2	C	1005	MET
2	C	1008	LEU
2	C	1009	ASP
2	C	1017	ASP
1	B	157	SER
1	B	161	GLU
1	B	297	ASP
1	B	306	ARG
1	B	340	LYS
1	B	375	VAL
1	B	396	ARG
1	B	598	LEU
1	B	611	TYR
2	D	1009	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	492	ASN

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)

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 (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	652/705 (92%)	-0.22	24 (3%) 41 39	19, 35, 68, 107	0
1	B	652/705 (92%)	-0.45	7 (1%) 80 79	15, 25, 52, 101	0
2	C	17/24 (70%)	0.21	0 100 100	38, 47, 71, 75	0
2	D	18/24 (75%)	-0.18	0 100 100	33, 42, 62, 66	0
All	All	1339/1458 (91%)	-0.32	31 (2%) 60 58	15, 30, 65, 107	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	686	GLN	5.4
1	A	686	GLN	4.6
1	A	395	SER	4.4
1	B	290	GLU	4.4
1	A	276	GLN	4.1
1	A	273	THR	4.0
1	A	278	ILE	3.8
1	A	478	ARG	3.7
1	B	273	THR	3.7
1	A	426	GLN	3.6
1	A	275	PRO	3.5
1	A	480	THR	3.4
1	A	427	LEU	3.1
1	B	250	ALA	2.9
1	A	396	ARG	2.9
1	A	98	ASP	2.7
1	A	393	ASP	2.6
1	A	547	VAL	2.5
1	A	669	PRO	2.5
1	A	476	TYR	2.5
1	A	482	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	276	GLN	2.4
1	A	297[A]	ASP	2.4
1	B	254	GLU	2.3
1	A	159	LYS	2.3
1	B	291	GLY	2.2
1	A	237	LYS	2.2
1	A	479	PHE	2.1
1	A	667	PRO	2.1
1	A	394	LEU	2.0
1	A	481	SER	2.0

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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.