



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 05:25 am BST

PDB ID : 2PV2
Title : Crystallographic Structure of SurA first peptidyl-prolyl isomerase domain complexed with peptide NFTLKFWDFRK
Authors : Xu, X.; McKay, D.B.
Deposited on : 2007-05-09
Resolution : 1.30 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

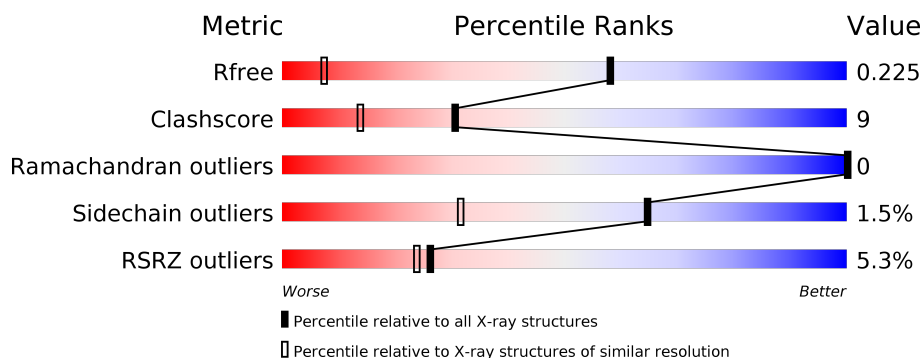
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.30 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	1058 (1.30-1.30)
Clashscore	141614	1101 (1.30-1.30)
Ramachandran outliers	138981	1058 (1.30-1.30)
Sidechain outliers	138945	1058 (1.30-1.30)
RSRZ outliers	127900	1029 (1.30-1.30)

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 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	103	<div> <div>3%</div> <div> <div></div> <div>88%</div> <div>12%</div> </div> </div>
1	B	103	<div> <div>4%</div> <div> <div></div> <div>81%</div> <div>19%</div> </div> </div>
1	C	103	<div> <div>3%</div> <div> <div></div> <div>89%</div> <div>11%</div> </div> </div>
1	D	103	<div> <div>12%</div> <div> <div></div> <div>76%</div> <div>23%</div> <div>.</div> </div> </div>
2	E	12	<div> <div></div> <div> <div></div> <div>83%</div> <div>17%</div> </div> </div>
2	F	12	<div> <div>8%</div> <div> <div></div> <div>83%</div> <div>17%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3722 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 Chaperone surA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	103	Total	C	N	O	S	0	0	0
			773	480	143	149	1			
1	B	103	Total	C	N	O	S	0	0	0
			773	480	143	149	1			
1	C	103	Total	C	N	O	S	0	0	0
			773	480	143	149	1			
1	D	103	Total	C	N	O	S	0	0	0
			773	480	143	149	1			

- Molecule 2 is a protein called C-peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	12	Total	C	N	O	0	0	0
			116	80	19	17			
2	F	12	Total	C	N	O	0	0	0
			116	80	19	17			

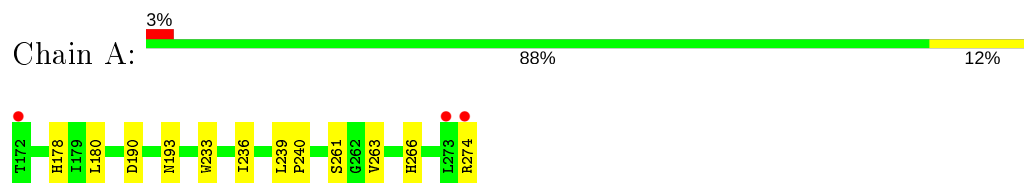
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	106	Total	O	0	0
			106	106		
3	B	96	Total	O	0	0
			96	96		
3	C	85	Total	O	0	0
			85	85		
3	D	89	Total	O	0	0
			89	89		
3	E	10	Total	O	0	0
			10	10		
3	F	12	Total	O	0	0
			12	12		

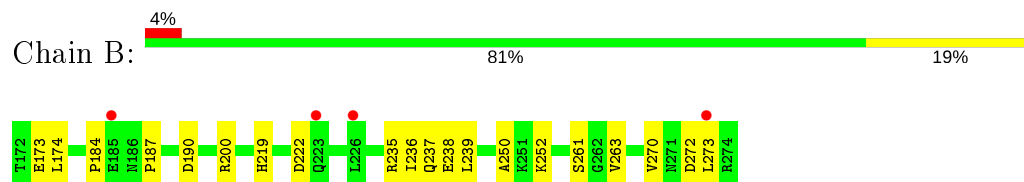
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

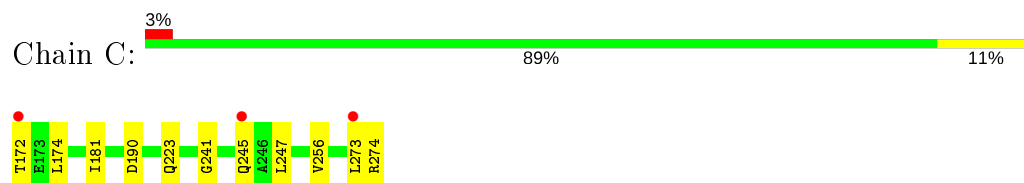
- Molecule 1: Chaperone surA



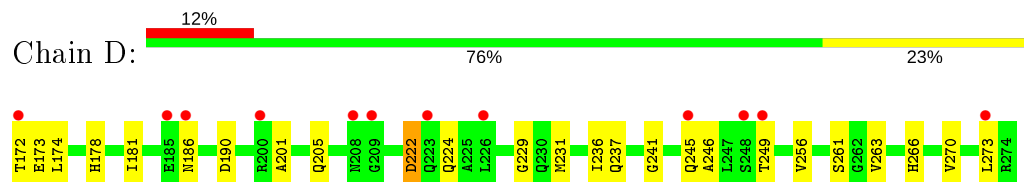
- Molecule 1: Chaperone surA



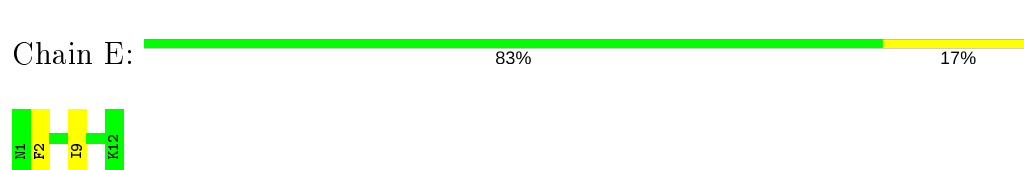
- Molecule 1: Chaperone surA




- Molecule 1: Chaperone surA

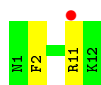


- Molecule 2: C-peptide



- Molecule 2: C-peptide

Chain F:  8% 83% 17%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.46 Å 50.63 Å 74.06 Å 90.00° 111.54° 90.00°	Depositor
Resolution (Å)	28.20 – 1.30 28.23 – 1.29	Depositor EDS
% Data completeness (in resolution range)	94.4 (28.20-1.30) 93.7 (28.23-1.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.89 (at 1.30 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.208 , 0.227 0.207 , 0.225	Depositor DCC
R_{free} test set	9256 reflections (9.93%)	wwPDB-VP
Wilson B-factor (Å ²)	14.1	Xtriage
Anisotropy	0.444	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 45.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.022 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3722	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 6.57% 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	0.34	0/785	0.72	1/1061 (0.1%)
1	B	0.32	0/785	0.69	1/1061 (0.1%)
1	C	0.31	0/785	0.70	2/1061 (0.2%)
1	D	0.29	0/785	0.69	2/1061 (0.2%)
2	E	0.47	0/120	0.60	0/158
2	F	0.44	0/120	0.60	0/158
All	All	0.33	0/3380	0.69	6/4560 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	190	ASP	CB-CG-OD2	5.23	123.01	118.30
1	C	190	ASP	CB-CG-OD2	5.19	122.97	118.30
1	B	190	ASP	CB-CG-OD2	5.16	122.94	118.30
1	D	190	ASP	CB-CG-OD2	5.13	122.92	118.30
1	C	181	ILE	N-CA-C	-5.13	97.16	111.00
1	D	181	ILE	N-CA-C	-5.08	97.29	111.00

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	773	0	770	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	773	0	770	21	0
1	C	773	0	770	8	0
1	D	773	0	770	23	0
2	E	116	0	117	1	0
2	F	116	0	117	1	0
3	A	106	0	0	2	0
3	B	96	0	0	2	0
3	C	85	0	0	2	0
3	D	89	0	0	1	0
3	E	10	0	0	0	0
3	F	12	0	0	0	0
All	All	3722	0	3314	61	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 9.

All (61) 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:174:LEU:HD22	1:B:273:LEU:HD22	1.16	1.15
1:D:174:LEU:HD22	1:D:273:LEU:HD22	1.35	1.03
1:B:174:LEU:HD22	1:B:273:LEU:CD2	2.02	0.89
1:D:174:LEU:HD22	1:D:273:LEU:CD2	2.08	0.83
1:A:236:ILE:HD12	1:A:239:LEU:HD12	1.64	0.79
1:B:174:LEU:CD2	1:B:273:LEU:HD22	2.07	0.79
1:B:270:VAL:CG1	1:B:273:LEU:HD21	2.16	0.76
1:D:178:HIS:HB3	1:D:231:MET:HE3	1.67	0.75
1:D:178:HIS:NE2	1:D:266:HIS:HD2	1.85	0.73
1:B:236:ILE:HD12	1:B:239:LEU:HD12	1.70	0.73
1:A:178:HIS:NE2	1:A:266:HIS:HD2	1.86	0.72
1:A:180:LEU:HD21	3:A:379:HOH:O	1.89	0.72
1:A:261:SER:OG	1:A:266:HIS:HE1	1.74	0.70
1:C:174:LEU:CD2	1:C:273:LEU:HD13	2.23	0.68
1:C:174:LEU:HD23	1:C:273:LEU:HD13	1.76	0.67
1:D:261:SER:OG	1:D:266:HIS:HE1	1.78	0.66
1:B:270:VAL:HG11	1:B:273:LEU:HD21	1.79	0.64
1:D:270:VAL:CG1	1:D:273:LEU:HD21	2.28	0.64
1:D:270:VAL:HG11	1:D:273:LEU:HD21	1.78	0.64
1:D:172:THR:HB	1:D:236:ILE:HD12	1.81	0.63
1:B:270:VAL:HG13	1:B:273:LEU:HD21	1.81	0.61
1:D:186:ASN:HB2	3:D:344:HOH:O	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:174:LEU:CD2	1:D:273:LEU:HD22	2.22	0.59
1:D:249:THR:O	1:D:249:THR:HG22	2.02	0.59
1:A:263:VAL:HB	3:A:379:HOH:O	2.03	0.58
1:B:235:ARG:HE	1:B:238:GLU:CD	2.07	0.57
1:A:233:TRP:CZ3	1:A:274:ARG:HG2	2.41	0.55
1:B:272:ASP:O	1:B:273:LEU:HD23	2.06	0.55
1:B:219:HIS:HE1	3:B:303:HOH:O	1.89	0.55
1:D:178:HIS:HB3	1:D:231:MET:CE	2.36	0.53
1:D:241:GLY:O	1:D:245:GLN:HG3	2.08	0.53
1:B:200:ARG:HG2	1:B:200:ARG:HH11	1.76	0.50
1:B:173:GLU:O	1:B:174:LEU:HD23	2.12	0.49
1:D:222:ASP:OD2	1:D:222:ASP:C	2.51	0.49
1:B:252:LYS:HE2	1:B:272:ASP:HA	1.94	0.49
1:B:184:PRO:O	1:B:187:PRO:HD3	2.13	0.48
1:D:173:GLU:O	1:D:174:LEU:HD23	2.13	0.48
1:D:224:GLN:HB2	1:D:231:MET:HE2	1.96	0.48
1:B:250:ALA:HB1	1:B:270:VAL:HG21	1.95	0.48
1:D:174:LEU:HD21	1:D:236:ILE:HD11	1.96	0.47
1:C:172:THR:HG22	3:C:297:HOH:O	2.14	0.46
1:C:273:LEU:HD12	1:C:274:ARG:N	2.30	0.46
1:A:178:HIS:NE2	1:A:266:HIS:CD2	2.77	0.46
2:F:11:ARG:HH11	2:F:11:ARG:HG3	1.80	0.45
1:D:201:ALA:O	1:D:205:GLN:HG3	2.16	0.45
1:D:246:ALA:HB1	1:D:256:VAL:HG11	1.98	0.45
1:C:174:LEU:HD21	1:C:273:LEU:HD13	1.96	0.44
1:C:223:GLN:HG2	3:C:344:HOH:O	2.17	0.44
1:C:241:GLY:O	1:C:245:GLN:HG3	2.18	0.44
1:B:200:ARG:HG2	1:B:200:ARG:NH1	2.32	0.44
1:B:270:VAL:HG13	1:B:273:LEU:CD2	2.47	0.43
1:C:247:LEU:HD23	1:C:256:VAL:HG21	2.00	0.43
1:B:261:SER:OG	1:B:263:VAL:HG22	2.19	0.42
1:D:237:GLN:CD	1:D:237:GLN:H	2.23	0.42
1:B:270:VAL:CG1	1:B:273:LEU:CD2	2.93	0.42
1:D:229:GLY:O	1:D:231:MET:CE	2.68	0.42
1:A:178:HIS:CE1	2:E:9:ILE:HD13	2.56	0.41
1:D:229:GLY:O	1:D:231:MET:HE3	2.19	0.41
1:D:261:SER:OG	1:D:263:VAL:HG22	2.20	0.41
1:A:240:PRO:HA	1:B:239:LEU:O	2.21	0.41
1:B:237:GLN:HG2	3:B:298:HOH:O	2.20	0.40

There are no symmetry-related clashes.

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	101/103 (98%)	101 (100%)	0	0	100	100
1	B	101/103 (98%)	100 (99%)	1 (1%)	0	100	100
1	C	101/103 (98%)	101 (100%)	0	0	100	100
1	D	101/103 (98%)	101 (100%)	0	0	100	100
2	E	10/12 (83%)	10 (100%)	0	0	100	100
2	F	10/12 (83%)	10 (100%)	0	0	100	100
All	All	424/436 (97%)	423 (100%)	1 (0%)	0	100	100

There are no Ramachandran outliers to report.

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	80/80 (100%)	79 (99%)	1 (1%)	69	35
1	B	80/80 (100%)	79 (99%)	1 (1%)	69	35
1	C	80/80 (100%)	80 (100%)	0	100	100
1	D	80/80 (100%)	79 (99%)	1 (1%)	69	35
2	E	12/12 (100%)	11 (92%)	1 (8%)	11	0
2	F	12/12 (100%)	11 (92%)	1 (8%)	11	0
All	All	344/344 (100%)	339 (98%)	5 (2%)	65	31

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	193	ASN
1	B	222	ASP
1	D	222	ASP
2	E	2	PHE
2	F	2	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	193	ASN
1	A	230	GLN
1	A	266	HIS
1	A	271	ASN
1	B	205	GLN
1	B	208	ASN
1	B	219	HIS
1	B	245	GLN
1	C	271	ASN
1	D	224	GLN
1	D	266	HIS
1	D	271	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 ⓘ

There are no ligands in this entry.

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, 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	103/103 (100%)	0.35	3 (2%) 51 49	9, 12, 21, 29	0
1	B	103/103 (100%)	0.50	4 (3%) 39 36	10, 16, 23, 32	0
1	C	103/103 (100%)	0.41	3 (2%) 51 49	11, 16, 27, 32	0
1	D	103/103 (100%)	0.69	12 (11%) 4 3	11, 18, 28, 33	0
2	E	12/12 (100%)	0.58	0 100 100	11, 14, 29, 29	0
2	F	12/12 (100%)	0.81	1 (8%) 11 8	13, 16, 33, 33	0
All	All	436/436 (100%)	0.50	23 (5%) 26 24	9, 16, 27, 33	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	185	GLU	6.3
1	B	226	LEU	4.7
1	B	185	GLU	3.6
1	C	273	LEU	3.6
1	D	248	SER	3.5
1	A	172	THR	3.2
1	D	249	THR	3.2
1	C	245	GLN	3.0
1	D	172	THR	2.9
1	D	208	ASN	2.8
1	B	273	LEU	2.5
1	D	226	LEU	2.4
2	F	11	ARG	2.4
1	D	245	GLN	2.4
1	B	223	GLN	2.4
1	D	209	GLY	2.4
1	D	200	ARG	2.4
1	D	186	ASN	2.2
1	A	274	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	172	THR	2.1
1	A	273	LEU	2.1
1	D	273	LEU	2.1
1	D	223	GLN	2.1

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.