



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 05:53 AM UTC

PDB ID : 3SZE / pdb_00003sze
Title : Crystal structure of the passenger domain of the E. coli autotransporter EspP
Authors : Khan, S.; Mian, H.S.; Sandercock, L.E.; Battaile, K.P.; Lam, R.; Chirgadze, N.Y.; Pai, E.F.
Deposited on : 2011-07-18
Resolution : 2.50 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	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

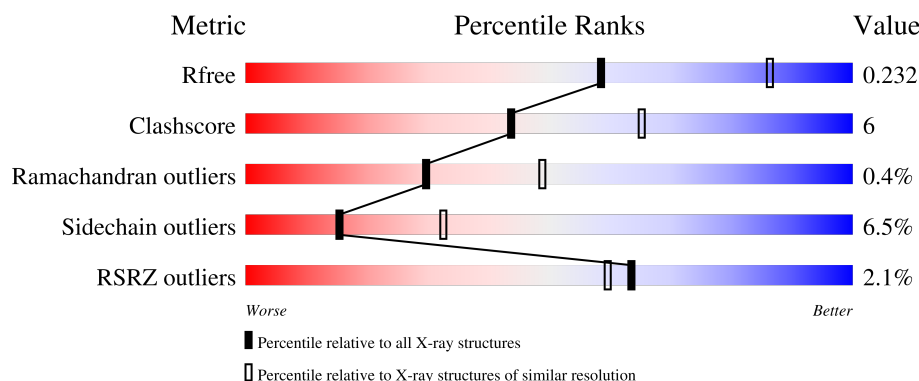
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

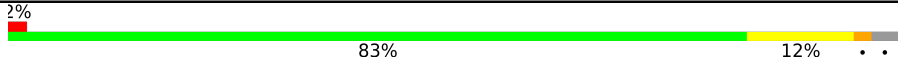
The reported resolution of this entry is 2.50 Å.

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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	968	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7576 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 Serine protease espP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	943	Total	C	N	O	S	0	1	0
			7252	4512	1251	1477	12			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	263	ALA	SER	engineered mutation	UNP Q7BSW5

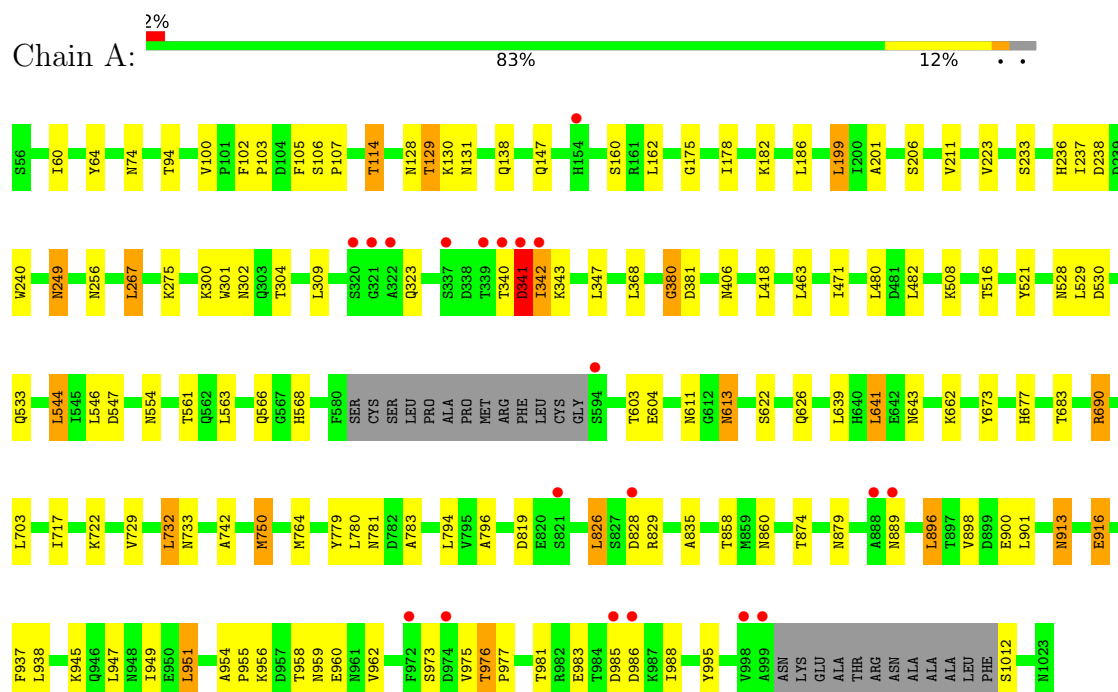
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	324	Total	O	0	0
			324	324		

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: Serine protease espP



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	88.34Å 88.34Å 311.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.50 30.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (30.00-2.50) 99.7 (30.00-2.50)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.47 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.188 , 0.231 0.192 , 0.232	Depositor DCC
R_{free} test set	2193 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	30.2	Xtriage
Anisotropy	0.182	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 34.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7576	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% 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

5.1 Standard geometry

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	0.51	0/7383	0.69	1/9998 (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	223	VAL	N-CA-C	8.04	114.81	106.21

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	7252	0	6999	87	0
2	A	324	0	0	3	0
All	All	7576	0	6999	87	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 6.

All (87) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance ($\text{\AA}$)	Clash overlap ($\text{\AA}$)
1:A:342:ILE:HD13	1:A:347:LEU:HD21	1.37	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:ASP:HA	1:A:342:ILE:CG2	1.95	0.97
1:A:114:THR:HG21	1:A:138:GLN:NE2	1.80	0.96
1:A:74:ASN:HD21	1:A:105:PHE:H	1.25	0.85
1:A:342:ILE:CD1	1:A:347:LEU:HD21	2.08	0.83
1:A:341:ASP:HA	1:A:342:ILE:HG22	1.61	0.79
1:A:114:THR:HG21	1:A:138:GLN:HE21	1.47	0.78
1:A:983:GLU:HG2	1:A:988:ILE:HG22	1.66	0.77
1:A:613:ASN:N	1:A:613:ASN:HD22	1.83	0.77
1:A:613:ASN:HD22	1:A:613:ASN:H	1.35	0.75
1:A:341:ASP:CA	1:A:342:ILE:HG22	2.17	0.74
1:A:341:ASP:HA	1:A:342:ILE:HG23	1.68	0.72
1:A:114:THR:CG2	1:A:138:GLN:NE2	2.54	0.71
1:A:201:ALA:HB1	1:A:267:LEU:HD22	1.72	0.69
1:A:732:LEU:HD22	1:A:733:ASN:OD1	1.92	0.69
1:A:528:ASN:HD22	1:A:554:ASN:HD21	1.43	0.65
1:A:566:GLN:HE22	1:A:673:TYR:H	1.44	0.63
1:A:147:GLN:NE2	1:A:160:SER:OG	2.31	0.63
1:A:60:ILE:HD11	1:A:603:THR:HG22	1.81	0.62
1:A:342:ILE:HD13	1:A:347:LEU:CD2	2.23	0.61
1:A:764:MET:HE2	1:A:794:LEU:HD11	1.84	0.59
1:A:182:LYS:HZ2	1:A:240:TRP:H	1.50	0.57
1:A:954:ALA:HB1	1:A:955:PRO:HD2	1.86	0.57
1:A:703:LEU:HD23	1:A:722:LYS:HB2	1.87	0.56
1:A:896:LEU:HD22	1:A:898:VAL:HG23	1.87	0.56
1:A:249:ASN:HD22	1:A:249:ASN:N	2.04	0.55
1:A:199:LEU:HD13	1:A:237:ILE:HB	1.89	0.55
1:A:860:ASN:HD21	1:A:879:ASN:HD22	1.55	0.54
1:A:60:ILE:HD11	1:A:603:THR:CG2	2.38	0.54
1:A:340:THR:O	1:A:341:ASP:HB3	2.06	0.54
1:A:508:LYS:HD3	1:A:530:ASP:OD2	2.08	0.54
1:A:913:ASN:HD21	1:A:916:GLU:H	1.56	0.53
1:A:956:LYS:HA	1:A:988:ILE:HG23	1.90	0.53
1:A:463:LEU:HD13	1:A:471:ILE:HD11	1.91	0.52
1:A:211:VAL:HG21	1:A:603:THR:HG21	1.91	0.52
1:A:147:GLN:HE21	1:A:160:SER:CB	2.22	0.52
1:A:983:GLU:CG	1:A:988:ILE:HG22	2.38	0.51
1:A:566:GLN:NE2	1:A:673:TYR:H	2.08	0.50
1:A:463:LEU:CD1	1:A:471:ILE:HD11	2.41	0.50
1:A:781:ASN:ND2	1:A:783:ALA:H	2.09	0.50
1:A:233:SER:HB3	2:A:1414:HOH:O	2.11	0.49
1:A:302:ASN:HD22	1:A:302:ASN:C	2.20	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:ALA:HB1	1:A:267:LEU:CD2	2.42	0.49
1:A:717:ILE:HG22	1:A:742:ALA:HB1	1.94	0.49
1:A:74:ASN:ND2	1:A:105:PHE:H	2.04	0.48
1:A:547:ASP:OD1	1:A:568:HIS:HD2	1.97	0.47
1:A:566:GLN:HE22	1:A:673:TYR:N	2.10	0.47
1:A:342:ILE:C	1:A:342:ILE:HD12	2.39	0.47
1:A:233:SER:HB2	1:A:256:ASN:HB3	1.97	0.47
1:A:544:LEU:HD13	1:A:563:LEU:CD2	2.44	0.47
1:A:236:HIS:HD2	1:A:238:ASP:OD2	1.98	0.46
1:A:764:MET:CE	1:A:794:LEU:HD11	2.44	0.46
1:A:732:LEU:HD22	1:A:733:ASN:CG	2.41	0.45
1:A:976:THR:HG23	1:A:977:PRO:HD2	1.97	0.45
1:A:129:THR:HG22	1:A:131:ASN:H	1.81	0.45
1:A:913:ASN:C	1:A:913:ASN:HD22	2.25	0.45
1:A:764:MET:O	1:A:796:ALA:HA	2.16	0.45
1:A:106:SER:N	1:A:107:PRO:CD	2.81	0.44
1:A:613:ASN:N	1:A:613:ASN:ND2	2.55	0.44
1:A:639:LEU:HG	1:A:641:LEU:HD13	2.00	0.43
1:A:826:LEU:HD21	1:A:835:ALA:HB1	2.00	0.43
1:A:102:PHE:CG	1:A:103:PRO:HD2	2.53	0.43
1:A:975:VAL:HG13	1:A:995:TYR:HB2	2.00	0.43
1:A:643:ASN:HA	1:A:662:LYS:O	2.18	0.43
1:A:683:THR:CG2	1:A:690:ARG:HD2	2.49	0.43
1:A:521:TYR:HA	2:A:1105:HOH:O	2.19	0.43
1:A:781:ASN:HD22	1:A:783:ALA:H	1.66	0.43
1:A:677:HIS:CD2	1:A:690:ARG:HG3	2.53	0.42
1:A:937:PHE:CE1	1:A:975:VAL:HG12	2.53	0.42
1:A:611:ASN:HB2	1:A:613:ASN:ND2	2.35	0.42
1:A:729:VAL:HB	1:A:750:MET:HG2	2.02	0.42
1:A:128:ASN:O	1:A:129:THR:CB	2.68	0.41
1:A:533:GLN:NE2	1:A:561:THR:OG1	2.53	0.41
1:A:626:GLN:NE2	2:A:1230:HOH:O	2.52	0.41
1:A:958:THR:HB	1:A:988:ILE:HD11	2.03	0.41
1:A:341:ASP:N	1:A:342:ILE:HG22	2.35	0.41
1:A:949:ILE:O	1:A:951:LEU:HD13	2.21	0.41
1:A:874:THR:HG22	1:A:900:GLU:HB2	2.01	0.41
1:A:129:THR:HG22	1:A:130:LYS:N	2.36	0.41
1:A:64:TYR:HA	1:A:604:GLU:OE2	2.21	0.41
1:A:175:GLY:HA3	1:A:301:TRP:CE2	2.56	0.41
1:A:178:ILE:HG21	1:A:300:LYS:HB3	2.03	0.41
1:A:622:SER:HA	1:A:626:GLN:HE22	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:GLY:O	1:A:381:ASP:C	2.64	0.40
1:A:683:THR:HG22	1:A:690:ARG:HD2	2.03	0.40
1:A:959:ASN:O	1:A:962:VAL:HG23	2.21	0.40
1:A:750:MET:HB2	1:A:779:TYR:HD2	1.87	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	938/968 (97%)	887 (95%)	47 (5%)	4 (0%)	30	49

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	341	ASP
1	A	129	THR
1	A	323	GLN
1	A	380	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	791/810 (98%)	740 (94%)	51 (6%)	16	33

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94	THR
1	A	100	VAL
1	A	114	THR
1	A	162	LEU
1	A	186	LEU
1	A	199	LEU
1	A	206	SER
1	A	249	ASN
1	A	267	LEU
1	A	275	LYS
1	A	304	THR
1	A	309	LEU
1	A	341	ASP
1	A	342	ILE
1	A	343	LYS
1	A	368	LEU
1	A	406	ASN
1	A	418	LEU
1	A	480	LEU
1	A	482	LEU
1	A	516	THR
1	A	529	LEU
1	A	544	LEU
1	A	546	LEU
1	A	613	ASN
1	A	641	LEU
1	A	690	ARG
1	A	732	LEU
1	A	750	MET
1	A	780	LEU
1	A	819	ASP
1	A	826	LEU
1	A	828	ASP
1	A	829	ARG
1	A	858	THR
1	A	889	ASN
1	A	896	LEU
1	A	901	LEU
1	A	913	ASN
1	A	916	GLU
1	A	938	LEU
1	A	945	LYS

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Mol	Chain	Res	Type
1	A	947	LEU
1	A	951	LEU
1	A	960	GLU
1	A	973	SER
1	A	976	THR
1	A	981	THR
1	A	985	ASP
1	A	986	ASP
1	A	1012	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	ASN
1	A	138	GLN
1	A	147	GLN
1	A	151	ASN
1	A	221	ASN
1	A	236	HIS
1	A	249	ASN
1	A	250	GLN
1	A	288	ASN
1	A	302	ASN
1	A	394	ASN
1	A	401	ASN
1	A	406	ASN
1	A	445	ASN
1	A	515	ASN
1	A	524	ASN
1	A	528	ASN
1	A	533	GLN
1	A	541	ASN
1	A	560	ASN
1	A	566	GLN
1	A	568	HIS
1	A	613	ASN
1	A	626	GLN
1	A	653	ASN
1	A	691	GLN
1	A	697	ASN
1	A	699	GLN
1	A	740	ASN

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Mol	Chain	Res	Type
1	A	781	ASN
1	A	798	ASN
1	A	851	ASN
1	A	860	ASN
1	A	861	ASN
1	A	879	ASN
1	A	913	ASN
1	A	931	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 oligosaccharides 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	943/968 (97%)	-0.25	20 (2%) 63 59	17, 29, 55, 77	1 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	342	ILE	3.7
1	A	821	SER	3.4
1	A	322	ALA	3.3
1	A	889	ASN	3.0
1	A	972	PHE	2.7
1	A	154	HIS	2.6
1	A	986	ASP	2.5
1	A	998	VAL	2.5
1	A	341	ASP	2.4
1	A	888	ALA	2.3
1	A	999	ALA	2.3
1	A	974	ASP	2.3
1	A	337	SER	2.3
1	A	828	ASP	2.3
1	A	321	GLY	2.2
1	A	985	ASP	2.2
1	A	339	THR	2.1
1	A	320	SER	2.1
1	A	340	THR	2.1
1	A	594	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 oligosaccharides 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.