



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:04 AM UTC

PDB ID : 7SR4 / pdb_00007sr4
Title : Single chain trimer HLA-A*02:01 (H98L, Y108C) with HPV.16 E7 peptide
YMLDLQPETTDLYC
Authors : Finton, K.A.K.; Rupert, P.B.
Deposited on : 2021-11-07
Resolution : 2.59 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

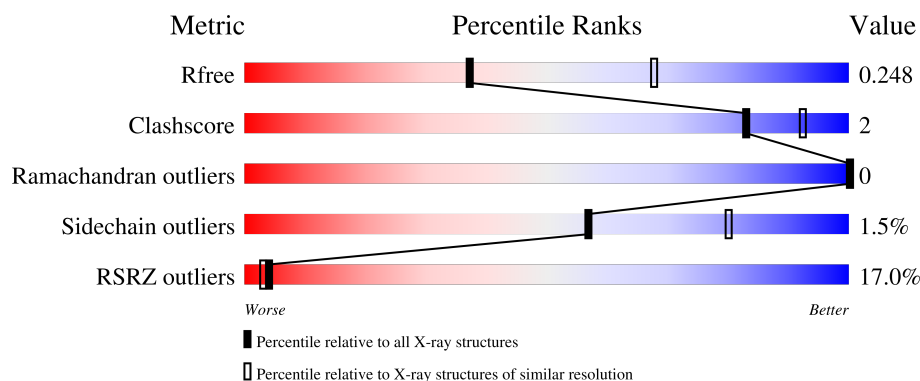
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (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	429	15% 73% 7% 20%
1	C	429	17% 83% 6% 11%
2	B	116	12% 91% 5%
2	D	116	10% 94%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7363 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 Protein E7 peptide,Beta-2-microglobulin,MHC class I antigen chimera.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	343	Total	C	N	O	S	0	0	0
			2657	1679	464	502	12			
1	C	382	Total	C	N	O	S	0	0	0
			2934	1851	517	553	13			

There are 86 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	14A	GLY	-	linker	UNP P03129
A	14B	CYS	-	linker	UNP P03129
A	14C	GLY	-	linker	UNP P03129
A	14D	GLY	-	linker	UNP P03129
A	14E	SER	-	linker	UNP P03129
A	14F	GLY	-	linker	UNP P03129
A	14G	GLY	-	linker	UNP P03129
A	14H	GLY	-	linker	UNP P03129
A	14I	GLY	-	linker	UNP P03129
A	14J	SER	-	linker	UNP P03129
A	14K	GLY	-	linker	UNP P03129
A	14L	GLY	-	linker	UNP P03129
A	14M	GLY	-	linker	UNP P03129
A	14N	GLY	-	linker	UNP P03129
A	14O	SER	-	linker	UNP P03129
A	124	GLY	-	linker	UNP P16213
A	125	GLY	-	linker	UNP P16213
A	126	GLY	-	linker	UNP P16213
A	127	GLY	-	linker	UNP P16213
A	128	SER	-	linker	UNP P16213
A	129	GLY	-	linker	UNP P16213
A	130	GLY	-	linker	UNP P16213
A	131	GLY	-	linker	UNP P16213
A	132	GLY	-	linker	UNP P16213

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Chain	Residue	Modelled	Actual	Comment	Reference
A	133	SER	-	linker	UNP P16213
A	134	GLY	-	linker	UNP P16213
A	135	GLY	-	linker	UNP P16213
A	136	GLY	-	linker	UNP P16213
A	137	GLY	-	linker	UNP P16213
A	138	SER	-	linker	UNP P16213
A	139	GLY	-	linker	UNP P16213
A	140	GLY	-	linker	UNP P16213
A	141	GLY	-	linker	UNP P16213
A	142	GLY	-	linker	UNP P16213
A	143	SER	-	linker	UNP P16213
A	217	LEU	HIS	engineered mutation	UNP A0A678ZGP6
A	227	CYS	TYR	engineered mutation	UNP A0A678ZGP6
A	419	HIS	-	expression tag	UNP A0A678ZGP6
A	420	HIS	-	expression tag	UNP A0A678ZGP6
A	421	HIS	-	expression tag	UNP A0A678ZGP6
A	422	HIS	-	expression tag	UNP A0A678ZGP6
A	423	HIS	-	expression tag	UNP A0A678ZGP6
A	424	HIS	-	expression tag	UNP A0A678ZGP6
C	14A	GLY	-	linker	UNP P03129
C	14B	CYS	-	linker	UNP P03129
C	14C	GLY	-	linker	UNP P03129
C	14D	GLY	-	linker	UNP P03129
C	14E	SER	-	linker	UNP P03129
C	14F	GLY	-	linker	UNP P03129
C	14G	GLY	-	linker	UNP P03129
C	14H	GLY	-	linker	UNP P03129
C	14I	GLY	-	linker	UNP P03129
C	14J	SER	-	linker	UNP P03129
C	14K	GLY	-	linker	UNP P03129
C	14L	GLY	-	linker	UNP P03129
C	14M	GLY	-	linker	UNP P03129
C	14N	GLY	-	linker	UNP P03129
C	14O	SER	-	linker	UNP P03129
C	124	GLY	-	linker	UNP P16213
C	125	GLY	-	linker	UNP P16213
C	126	GLY	-	linker	UNP P16213
C	127	GLY	-	linker	UNP P16213
C	128	SER	-	linker	UNP P16213
C	129	GLY	-	linker	UNP P16213
C	130	GLY	-	linker	UNP P16213
C	131	GLY	-	linker	UNP P16213

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Chain	Residue	Modelled	Actual	Comment	Reference
C	132	GLY	-	linker	UNP P16213
C	133	SER	-	linker	UNP P16213
C	134	GLY	-	linker	UNP P16213
C	135	GLY	-	linker	UNP P16213
C	136	GLY	-	linker	UNP P16213
C	137	GLY	-	linker	UNP P16213
C	138	SER	-	linker	UNP P16213
C	139	GLY	-	linker	UNP P16213
C	140	GLY	-	linker	UNP P16213
C	141	GLY	-	linker	UNP P16213
C	142	GLY	-	linker	UNP P16213
C	143	SER	-	linker	UNP P16213
C	217	LEU	HIS	engineered mutation	UNP A0A678ZGP6
C	227	CYS	TYR	engineered mutation	UNP A0A678ZGP6
C	419	HIS	-	expression tag	UNP A0A678ZGP6
C	420	HIS	-	expression tag	UNP A0A678ZGP6
C	421	HIS	-	expression tag	UNP A0A678ZGP6
C	422	HIS	-	expression tag	UNP A0A678ZGP6
C	423	HIS	-	expression tag	UNP A0A678ZGP6
C	424	HIS	-	expression tag	UNP A0A678ZGP6

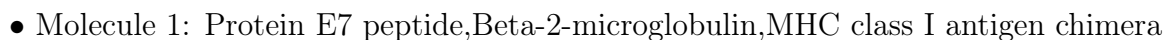
- Molecule 2 is a protein called VHH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	112	Total	C	N	O	S	0	0	0
			818	506	140	168	4			
2	D	111	Total	C	N	O	S	0	0	0
			812	502	139	167	4			

- Molecule 3 is water.

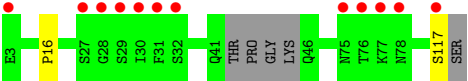
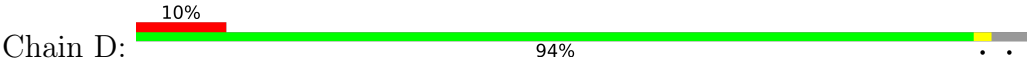
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	40	Total	O	0	0
			40	40		
3	B	22	Total	O	0	0
			22	22		
3	C	50	Total	O	0	0
			50	50		
3	D	30	Total	O	0	0
			30	30		

- Molecule 1: Protein E7 peptide, Beta-2-microglobulin, MHC class I antigen chimera





● Molecule 2: VHH



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	117.88Å 117.88Å 262.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.91 – 2.59 48.91 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.91-2.59) 99.8 (48.91-2.59)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 2.58Å)	Xtriage
Refinement program	PHENIX 1.19.1_4122	Depositor
R, R_{free}	0.218 , 0.247 0.220 , 0.248	Depositor DCC
R_{free} test set	2962 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	46.5	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7363	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 23.23 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.8612e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.11	0/2728	0.29	0/3714
1	C	0.13	0/3016	0.30	0/4116
2	B	0.11	0/831	0.32	0/1126
2	D	0.11	0/825	0.30	0/1119
All	All	0.12	0/7400	0.30	0/10075

There are no bond length outliers.

There are no bond angle outliers.

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	2657	0	2382	15	0
1	C	2934	0	2600	15	0
2	B	818	0	749	2	0
2	D	812	0	739	1	0
3	A	40	0	0	0	0
3	B	22	0	0	0	0
3	C	50	0	0	0	0
3	D	30	0	0	0	0
All	All	7363	0	6470	33	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 2.

All (33) 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:B:68:ARG:NH2	2:B:91:ASP:OD2	2.24	0.69
1:A:276:TRP:HB2	1:A:287:LYS:HG3	1.85	0.58
1:C:300:ARG:NH1	1:C:304:GLU:OE2	2.37	0.58
1:A:2:MET:HE1	1:A:210:VAL:HB	1.85	0.57
1:C:149:ARG:NH2	1:C:245:ASP:OD1	2.38	0.57
1:A:358:LEU:HD12	1:A:386:LYS:HG2	1.87	0.56
1:A:178:ARG:C	1:A:188:MET:HE3	2.35	0.52
1:A:77:ASP:OD2	1:A:178:ARG:HD2	2.11	0.50
1:C:308:VAL:O	1:C:312:ARG:HG3	2.12	0.50
2:B:41:GLN:NE2	2:B:45:LYS:O	2.44	0.50
1:A:308:VAL:O	1:A:312:ARG:HG3	2.12	0.48
1:C:276:TRP:HB2	1:C:287:LYS:HD2	1.96	0.48
1:C:245:ASP:OD2	1:C:256:TYR:OH	2.23	0.46
1:A:171:VAL:HG11	1:A:322:LEU:HD13	1.97	0.46
1:A:224:LEU:HD13	1:A:261:TYR:CD1	2.52	0.45
1:A:6:GLN:O	1:A:9:THR:OG1	2.30	0.45
1:C:251:ARG:HG2	1:C:251:ARG:NH1	2.32	0.44
1:A:10:THR:O	1:A:10:THR:OG1	2.35	0.44
1:C:250:TRP:HB3	1:C:312:ARG:HD3	2.00	0.44
1:C:251:ARG:HG2	1:C:251:ARG:HH11	1.83	0.44
1:C:41:ASN:OD1	1:C:121:ARG:NH1	2.36	0.43
1:C:358:LEU:HD12	1:C:386:LYS:HD3	1.99	0.43
1:A:84:TRP:CE2	1:A:260:ALA:HB2	2.54	0.43
1:A:309:GLU:OE2	1:A:313:ARG:HD2	2.19	0.43
1:A:148:MET:HB2	1:A:311:LEU:HD13	2.00	0.42
1:C:2:MET:HE1	1:C:210:VAL:HB	2.01	0.42
1:C:224:LEU:HD13	1:C:261:TYR:CD1	2.55	0.41
1:A:250:TRP:HB3	1:A:312:ARG:HD3	2.03	0.41
1:C:252:PHE:CZ	1:C:304:GLU:HG2	2.55	0.41
2:D:16:PRO:HD3	2:D:117:SER:HB2	2.02	0.41
1:C:8:GLU:HA	1:C:11:ASP:HB2	2.02	0.41
1:C:84:TRP:CE2	1:C:260:ALA:HB2	2.56	0.40
1:A:10:THR:HG23	1:A:216:THR:HG21	2.04	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	331/429 (77%)	323 (98%)	8 (2%)	0	100	100
1	C	374/429 (87%)	365 (98%)	9 (2%)	0	100	100
2	B	108/116 (93%)	106 (98%)	2 (2%)	0	100	100
2	D	107/116 (92%)	106 (99%)	1 (1%)	0	100	100
All	All	920/1090 (84%)	900 (98%)	20 (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	261/353 (74%)	253 (97%)	8 (3%)	35	63
1	C	282/353 (80%)	281 (100%)	1 (0%)	84	93
2	B	84/97 (87%)	82 (98%)	2 (2%)	43	70
2	D	83/97 (86%)	83 (100%)	0	100	100
All	All	710/900 (79%)	699 (98%)	11 (2%)	57	80

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	THR
1	A	14	CYS

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Mol	Chain	Res	Type
1	A	154	SER
1	A	238	VAL
1	A	308	VAL
1	A	313	ARG
1	A	374	VAL
1	A	402	CYS
2	B	20	LEU
2	B	55	SER
1	C	402	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	HIS
2	B	85	ASN
1	C	37	HIS
1	C	175	GLN
1	C	197	GLN
1	C	230	GLN
1	C	331	HIS

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

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	343/429 (79%)	0.90	63 (18%) 3 3	34, 52, 101, 127	0
1	C	382/429 (89%)	1.09	72 (18%) 3 2	34, 54, 104, 134	0
2	B	112/116 (96%)	0.51	14 (12%) 8 6	32, 41, 86, 102	0
2	D	111/116 (95%)	0.50	12 (10%) 11 8	33, 41, 84, 100	0
All	All	948/1090 (86%)	0.88	161 (16%) 4 3	32, 49, 100, 134	0

All (161) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	340	HIS	9.6
2	B	75	ASN	9.1
2	D	75	ASN	7.4
1	C	14	CYS	7.1
1	C	417	TRP	7.1
1	C	25	ILE	6.8
1	A	410	PRO	6.8
1	C	342	ALA	6.7
1	A	347	TRP	6.5
1	C	122	ASP	6.2
2	B	3	GLU	6.1
1	C	13	TYR	5.9
1	C	335	HIS	5.8
1	C	366	ASP	5.7
1	A	405	GLN	5.7
1	C	229	ASN	5.5
1	A	331	HIS	5.5
2	D	29	SER	5.5
1	A	404	VAL	5.5
1	A	25	ILE	5.5
1	C	341	GLU	5.3

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Mol	Chain	Res	Type	RSRZ
1	A	373	LEU	5.2
1	C	144	GLY	5.2
1	C	413	LEU	5.1
2	D	3	GLU	5.1
1	C	7	PRO	5.0
2	B	29	SER	5.0
1	A	14	CYS	4.9
1	C	370	ASP	4.9
1	C	398	GLN	4.9
1	C	363	ASP	4.8
2	B	76	THR	4.8
1	A	13	TYR	4.6
1	C	360	TRP	4.5
1	C	407	GLU	4.5
2	B	31	PHE	4.5
1	C	368	THR	4.5
2	D	30	ILE	4.4
1	A	401	THR	4.3
2	B	28	GLY	4.3
1	A	144	GLY	4.3
2	B	77	LYS	4.2
1	C	397	GLU	4.2
1	A	328	PRO	4.2
1	A	402	CYS	4.2
1	A	387	TRP	4.2
1	C	367	GLN	4.2
1	A	372	GLU	4.1
1	C	281	MET	4.1
1	A	407	GLU	4.1
1	A	160	ARG	4.0
2	B	78	ASN	4.0
2	B	30	ILE	4.0
1	C	6	GLN	4.0
1	C	160	ARG	3.9
1	A	8	GLU	3.9
2	D	76	THR	3.9
1	C	332	MET	3.8
1	A	409	LEU	3.8
1	C	347	TRP	3.8
1	C	331	HIS	3.8
1	A	11	ASP	3.8
1	A	390	VAL	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	396	GLN	3.7
1	C	373	LEU	3.7
1	A	358	LEU	3.7
1	A	10	THR	3.6
1	A	7	PRO	3.6
1	A	346	CYS	3.6
1	A	374	VAL	3.6
1	C	8	GLU	3.5
1	C	321	THR	3.5
1	A	121	ARG	3.4
1	A	6	GLN	3.3
2	D	31	PHE	3.3
1	A	357	THR	3.3
1	A	388	ALA	3.3
1	C	288	HIS	3.3
1	C	304	GLU	3.2
1	A	344	LEU	3.2
1	C	333	THR	3.2
1	A	99	LYS	3.1
2	D	117	SER	3.1
1	C	392	VAL	3.1
1	A	12	LEU	3.0
1	C	402	CYS	3.0
2	D	78	ASN	3.0
2	B	117	SER	3.0
1	A	355	GLU	2.9
1	C	387	TRP	2.9
1	C	12	LEU	2.9
1	C	371	THR	2.9
1	C	394	SER	2.9
1	C	279	ALA	2.9
1	A	197	GLN	2.9
1	A	256	TYR	2.9
1	C	256	TYR	2.9
1	C	346	CYS	2.8
1	C	369	GLN	2.8
1	A	345	ARG	2.8
2	D	27	SER	2.8
2	D	32	SER	2.8
1	C	415	LEU	2.8
1	C	343	THR	2.7
1	C	362	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	399	ARG	2.7
1	C	410	PRO	2.7
1	A	9	THR	2.7
1	A	233	ALA	2.7
1	C	10	THR	2.7
1	C	365	GLU	2.7
1	C	197	GLN	2.7
1	C	159	GLY	2.7
1	C	393	PRO	2.7
1	C	372	GLU	2.6
1	A	403	HIS	2.6
1	A	159	GLY	2.6
1	C	391	VAL	2.6
2	D	28	GLY	2.6
1	A	375	GLU	2.6
1	A	327	ALA	2.6
1	C	359	THR	2.6
1	A	98	GLU	2.5
1	A	184	ALA	2.5
1	A	325	THR	2.5
1	A	356	ILE	2.5
1	A	288	HIS	2.5
2	B	7	VAL	2.5
2	B	27	SER	2.4
1	C	400	TYR	2.4
1	A	281	MET	2.4
1	C	9	THR	2.4
1	C	329	LYS	2.4
1	C	162	GLU	2.4
1	C	259	TYR	2.3
1	A	279	ALA	2.3
1	C	327	ALA	2.3
1	A	161	GLY	2.3
1	A	330	THR	2.3
1	C	330	THR	2.3
2	B	32	SER	2.3
2	D	77	LYS	2.3
1	A	97	THR	2.3
1	A	406	HIS	2.3
1	A	229	ASN	2.2
1	C	297	GLU	2.2
1	A	232	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	271	GLU	2.2
1	A	408	GLY	2.2
1	C	344	LEU	2.2
1	A	41	ASN	2.1
2	B	45	LYS	2.1
1	A	43	LYS	2.1
1	C	158	PRO	2.1
1	A	42	GLY	2.1
1	C	240	ARG	2.0
1	C	408	GLY	2.0
1	A	149	ARG	2.0
1	C	121	ARG	2.0
1	A	234	GLY	2.0
1	A	182	ASP	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.