



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 27, 2016 – 05:07 AM EDT

PDB ID : 4ZVO  
Title : Caspase-7 Variant 4 (V4) with reprogrammed substrate specificity due to Y230V/W232Y/S234V/Q276D substitutions bound to VEID inhibitor.  
Authors : Hill, M.E.; MacPherson, D.J.; Hardy, J.A.  
Deposited on : 2015-05-18  
Resolution : 2.85 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/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  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20027790  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20027790

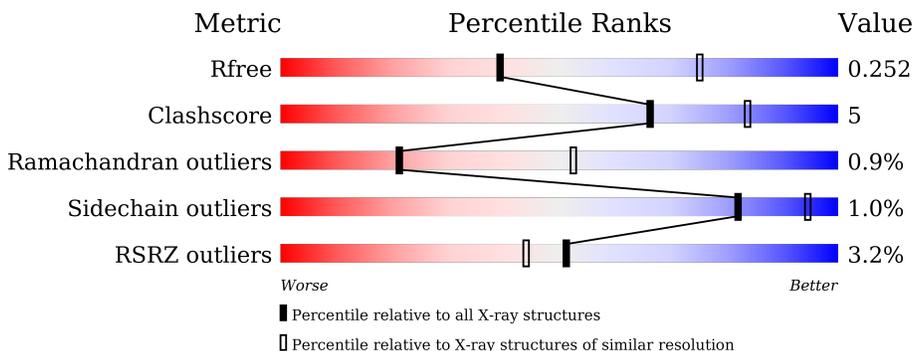
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.85 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))
$R_{free}$	91344	2228 (2.90-2.82)
Clashscore	102246	2499 (2.90-2.82)
Ramachandran outliers	100387	2439 (2.90-2.82)
Sidechain outliers	100360	2442 (2.90-2.82)
RSRZ outliers	91569	2236 (2.90-2.82)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	198	
1	C	198	
2	B	113	
2	D	113	
3	E	5	
3	F	5	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3792 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 Caspase-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	139	1092	687	187	207	11	0	0	0
1	C	140	1099	691	188	209	11	0	0	0

- Molecule 2 is a protein called Caspase-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	93	762	491	127	140	4	0	0	0
2	D	93	762	491	127	140	4	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	230	VAL	TYR	engineered mutation	UNP P55210
B	232	TYR	TRP	engineered mutation	UNP P55210
B	234	VAL	SER	engineered mutation	UNP P55210
B	276	ASP	GLN	engineered mutation	UNP P55210
B	304	LEU	-	expression tag	UNP P55210
B	305	GLU	-	expression tag	UNP P55210
B	306	HIS	-	expression tag	UNP P55210
B	307	HIS	-	expression tag	UNP P55210
B	308	HIS	-	expression tag	UNP P55210
B	309	HIS	-	expression tag	UNP P55210
B	310	HIS	-	expression tag	UNP P55210
B	311	HIS	-	expression tag	UNP P55210
D	530	VAL	TYR	engineered mutation	UNP P55210
D	532	TYR	TRP	engineered mutation	UNP P55210
D	534	VAL	SER	engineered mutation	UNP P55210
D	576	ASP	GLN	engineered mutation	UNP P55210

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
D	604	LEU	-	expression tag	UNP P55210
D	605	GLU	-	expression tag	UNP P55210
D	606	HIS	-	expression tag	UNP P55210
D	607	HIS	-	expression tag	UNP P55210
D	608	HIS	-	expression tag	UNP P55210
D	609	HIS	-	expression tag	UNP P55210
D	610	HIS	-	expression tag	UNP P55210
D	611	HIS	-	expression tag	UNP P55210

- Molecule 3 is a protein called Peptide ACE-VAL-GLU-ILE-ASJ.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace	
			Total	C	N				O
3	E	5	35	22	4	9	0	0	0
3	F	5	35	22	4	9	0	0	0

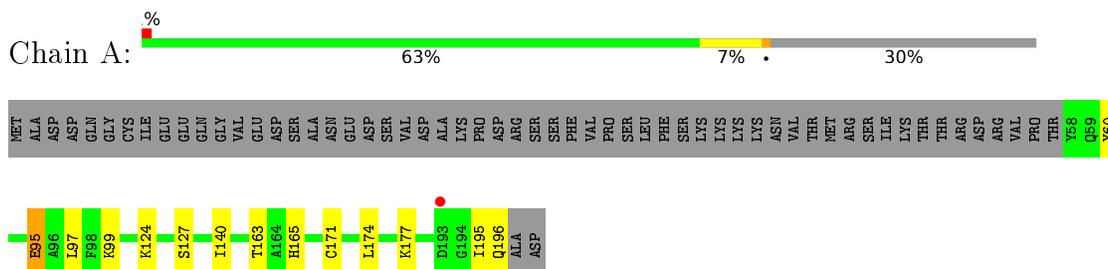
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total	O	0	0
			4	4		
4	B	1	Total	O	0	0
			1	1		
4	C	1	Total	O	0	0
			1	1		
4	D	1	Total	O	0	0
			1	1		

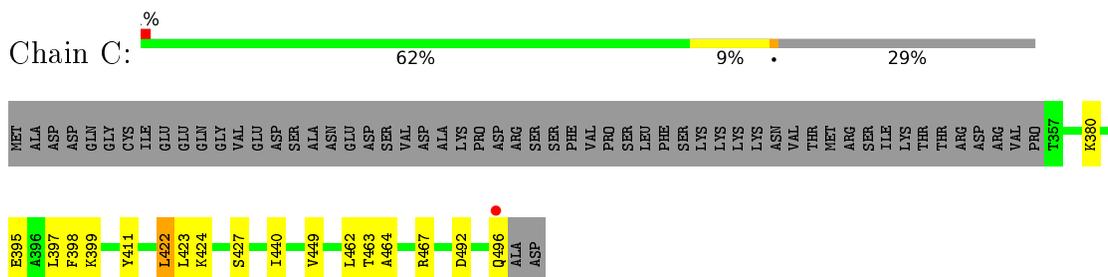
### 3 Residue-property plots [i](#)

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

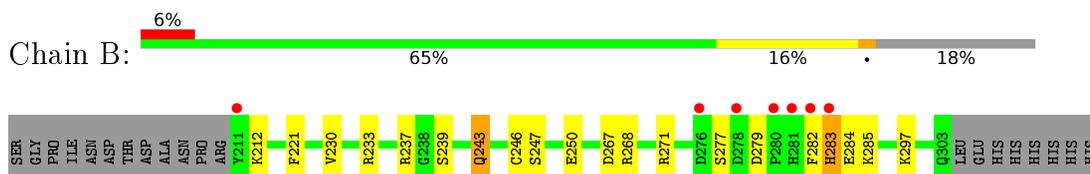
- Molecule 1: Caspase-7



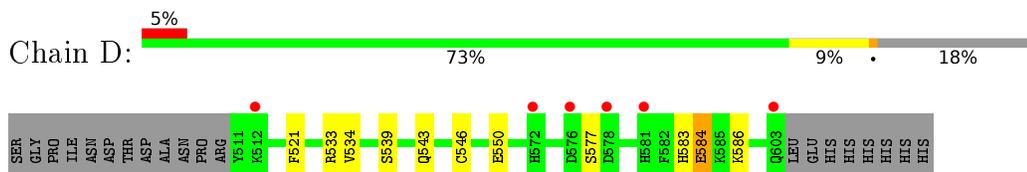
- Molecule 1: Caspase-7



- Molecule 2: Caspase-7



- Molecule 2: Caspase-7



- Molecule 3: Peptide ACE-VAL-GLU-ILE-ASJ



There are no outlier residues recorded for this chain.

- Molecule 3: Peptide ACE-VAL-GLU-ILE-ASJ

Chain F:  80% 20%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.27Å 88.27Å 187.28Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.22 – 2.85 38.22 – 2.85	Depositor EDS
% Data completeness (in resolution range)	98.5 (38.22-2.85) 98.5 (38.22-2.85)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.59 (at 2.85Å)	Xtrriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.209 , 0.251 0.211 , 0.252	Depositor DCC
$R_{free}$ test set	974 reflections (4.85%)	DCC
Wilson B-factor (Å <sup>2</sup> )	68.7	Xtrriage
Anisotropy	0.122	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 33.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.53$ , $\langle L^2 \rangle = 0.37$	Xtrriage
Estimated twinning fraction	0.001 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3792	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	65.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ASJ, ACE

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	0.29	0/1109	0.53	0/1486
1	C	0.30	0/1116	0.56	1/1496 (0.1%)
2	B	0.31	0/783	0.64	1/1060 (0.1%)
2	D	0.30	0/783	0.60	1/1060 (0.1%)
3	E	0.24	0/24	0.41	0/32
3	F	0.21	0/24	0.89	0/32
All	All	0.30	0/3839	0.58	3/5166 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	543	GLN	CA-CB-CG	-5.89	100.45	113.40
2	B	243	GLN	CA-CB-CG	5.53	125.56	113.40
1	C	422	LEU	CA-CB-CG	5.36	127.62	115.30

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	1092	0	1078	10	0
1	C	1099	0	1085	12	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	762	0	739	12	0
2	D	762	0	739	9	0
3	E	35	0	32	0	0
3	F	35	0	33	1	0
4	A	4	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
All	All	3792	0	3706	37	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 5.

All (37) 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:D:577:SER:HA	2:D:583:HIS:CD2	2.24	0.71
2:D:546:CYS:O	2:D:550:GLU:HG3	1.91	0.70
2:D:583:HIS:CD2	2:D:584:GLU:H	2.11	0.67
1:C:395:GLU:O	1:C:399:LYS:HG3	1.96	0.66
2:D:577:SER:HA	2:D:583:HIS:HD2	1.59	0.65
1:C:463:THR:HG21	2:D:521:PHE:HE2	1.63	0.63
1:C:423:LEU:HD12	1:C:462:LEU:HD22	1.82	0.60
1:A:174:LEU:HA	1:A:177:LYS:HD2	1.86	0.58
1:A:163:THR:HG21	2:B:221:PHE:HE1	1.70	0.57
1:A:195:ILE:O	1:A:196:GLN:HB3	2.06	0.56
2:B:282:PHE:O	2:B:285:LYS:HD3	2.07	0.54
2:D:533:ARG:HA	2:D:539:SER:HA	1.90	0.53
2:B:233:ARG:HA	2:B:239:SER:HA	1.92	0.52
1:A:97:LEU:HD13	1:A:140:ILE:HG21	1.92	0.51
2:B:246:CYS:O	2:B:250:GLU:HG3	2.10	0.51
1:C:397:LEU:HD13	1:C:440:ILE:HG21	1.92	0.51
1:C:464:ALA:O	1:C:467:ARG:HG3	2.10	0.51
2:B:247:SER:OG	2:B:268:ARG:NH1	2.45	0.50
1:A:60:TYR:CD2	2:B:297:LYS:HB2	2.48	0.48
1:A:163:THR:HG21	2:B:221:PHE:CE1	2.50	0.47
2:D:583:HIS:CD2	2:D:584:GLU:HG2	2.50	0.47
2:B:267:ASP:OD1	2:B:271:ARG:NH1	2.48	0.46
1:C:397:LEU:HD22	1:C:440:ILE:HD13	1.96	0.46
1:C:424:LYS:O	1:C:427:SER:OG	2.24	0.46
2:B:237:ARG:HB3	2:B:243:GLN:NE2	2.30	0.46

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:411:TYR:CG	1:C:422:LEU:HD11	2.51	0.46
2:B:212:LYS:HG3	1:C:496:GLN:O	2.17	0.45
1:A:95:GLU:O	1:A:99:LYS:HG3	2.16	0.45
2:B:279:ASP:HB3	2:B:283:HIS:H	1.81	0.45
2:D:534:VAL:HG22	3:F:1:VAL:HG12	1.99	0.45
1:A:124:LYS:O	1:A:127:SER:HB3	2.18	0.44
1:C:395:GLU:HA	1:C:398:PHE:CE2	2.53	0.43
1:A:127:SER:HB2	1:A:165:HIS:HB3	2.01	0.42
1:C:492:ASP:HB3	2:D:586:LYS:O	2.20	0.42
2:B:285:LYS:HA	2:B:285:LYS:HD2	1.56	0.41
1:C:411:TYR:CD1	1:C:422:LEU:HD11	2.55	0.41
1:A:171:CYS:SG	1:A:174:LEU:HD12	2.61	0.41

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	137/198 (69%)	132 (96%)	5 (4%)	0	100	100
1	C	138/198 (70%)	132 (96%)	6 (4%)	0	100	100
2	B	91/113 (80%)	84 (92%)	4 (4%)	3 (3%)	5	17
2	D	91/113 (80%)	85 (93%)	5 (6%)	1 (1%)	17	47
3	E	3/5 (60%)	3 (100%)	0	0	100	100
3	F	3/5 (60%)	3 (100%)	0	0	100	100
All	All	463/632 (73%)	439 (95%)	20 (4%)	4 (1%)	21	52

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	584	GLU

Continued on next page...

*Continued from previous page...*

Mol	Chain	Res	Type
2	B	277	SER
2	B	284	GLU
2	B	283	HIS

### 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	119/172 (69%)	118 (99%)	1 (1%)	86	96
1	C	120/172 (70%)	118 (98%)	2 (2%)	68	89
2	B	85/103 (82%)	84 (99%)	1 (1%)	78	93
2	D	85/103 (82%)	85 (100%)	0	100	100
3	E	3/3 (100%)	3 (100%)	0	100	100
3	F	3/3 (100%)	3 (100%)	0	100	100
All	All	415/556 (75%)	411 (99%)	4 (1%)	82	94

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	GLU
2	B	230	VAL
1	C	380	LYS
1	C	449	VAL

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

Mol	Chain	Res	Type
2	B	243	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
3	ASJ	E	4	1,3	4,7,7	2.53	1 (25%)	2,8,8	2.39	1 (50%)
3	ASJ	F	4	1,3	4,7,7	2.56	1 (25%)	2,8,8	2.78	1 (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
3	ASJ	E	4	1,3	-	0/4/6/6	0/0/0/0
3	ASJ	F	4	1,3	-	0/4/6/6	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	4	ASJ	O-C	-4.67	1.22	1.42
3	E	4	ASJ	O-C	-4.63	1.22	1.42

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	4	ASJ	O-C-CA	3.30	120.52	111.77
3	F	4	ASJ	O-C-CA	3.79	121.80	111.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	139/198 (70%)	-0.21	1 (0%) 89 88	51, 67, 82, 93	0
1	C	140/198 (70%)	-0.34	1 (0%) 89 88	46, 57, 73, 106	0
2	B	93/113 (82%)	0.25	7 (7%) 17 11	48, 62, 109, 118	0
2	D	93/113 (82%)	0.13	6 (6%) 22 16	46, 60, 106, 111	0
3	E	3/5 (60%)	-0.02	0 100 100	72, 72, 79, 79	0
3	F	3/5 (60%)	-0.64	0 100 100	60, 60, 67, 70	0
All	All	471/632 (74%)	-0.09	15 (3%) 51 44	46, 61, 95, 118	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	278	ASP	3.7
2	B	281	HIS	3.4
1	C	496	GLN	3.1
2	B	211	TYR	3.0
2	B	283	HIS	2.8
2	B	276	ASP	2.7
1	A	193	ASP	2.7
2	B	282	PHE	2.6
2	D	576	ASP	2.5
2	B	280	PRO	2.5
2	D	578	ASP	2.5
2	D	581	HIS	2.5
2	D	572	HIS	2.4
2	D	512	LYS	2.2
2	D	603	GLN	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	ASJ	E	4	8/8	0.97	0.13	-	64,68,70,70	0
3	ASJ	F	4	8/8	0.96	0.12	-	55,61,63,65	0

## 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.