



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 22, 2025 – 01:16 AM EDT

PDB ID : 6V8J / pdb\_00006v8j  
Title : Crystal structure of Ara h 8.0201  
Authors : Offermann, L.R.; Pote, S.; Hurlburt, B.K.; McBride, J.K.; Chruszcz, M.  
Deposited on : 2019-12-11  
Resolution : 1.95 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

<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 : 4.02b-467  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0rc1  
EDS : 3.0  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.006 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.42

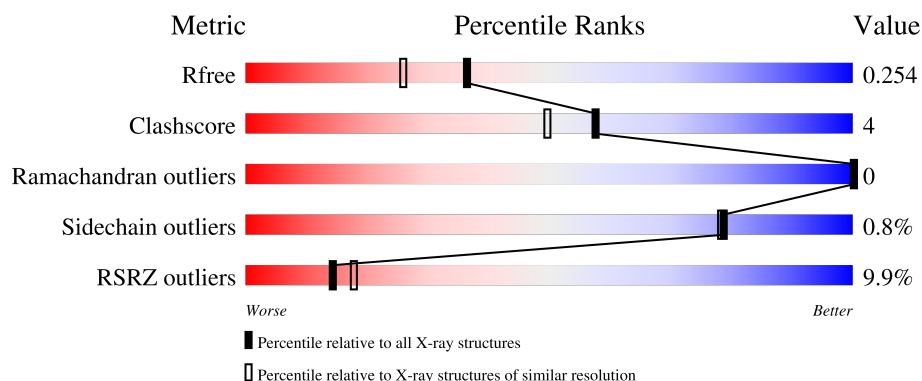
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.95 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))
$R_{free}$	164625	3187 (1.96-1.96)
Clashscore	180529	3412 (1.96-1.96)
Ramachandran outliers	177936	3390 (1.96-1.96)
Sidechain outliers	177891	3390 (1.96-1.96)
RSRZ outliers	164620	3186 (1.96-1.96)

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  $\geq 3$ , 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	153	<div> <div>13%</div> <div> <div></div> <div>87%</div> <div>12%</div> <div>.</div> </div> </div>
1	B	153	<div> <div>9%</div> <div> <div></div> <div>93%</div> <div>5%</div> <div>..</div> </div> </div>
1	C	153	<div> <div>10%</div> <div> <div></div> <div>87%</div> <div>12%</div> <div>.</div> </div> </div>
1	D	153	<div> <div>7%</div> <div> <div></div> <div>89%</div> <div>9%</div> <div>.</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4721 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 Ara h 8 allergen isoform.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	152	Total	C	N	O	0	0	0
			1133	720	182	231			
1	B	151	Total	C	N	O	0	0	0
			1127	716	181	230			
1	C	152	Total	C	N	O	0	0	0
			1133	718	182	233			
1	D	150	Total	C	N	O	0	0	0
			1131	720	179	232			

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		
2	B	1	Total	Na	0	0
			1	1		
2	C	1	Total	Na	0	0
			1	1		
2	D	1	Total	Na	0	0
			1	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

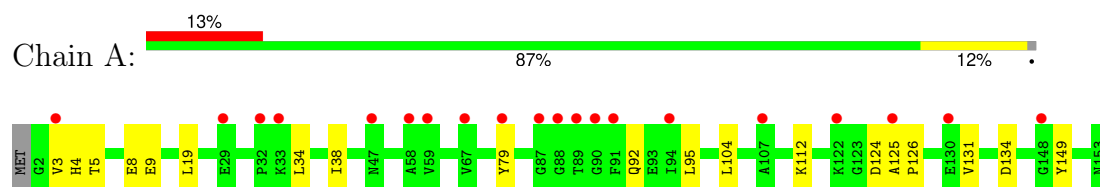
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	42	Total	O	0	0
			42	42		
4	B	41	Total	O	0	0
			41	41		
4	C	46	Total	O	0	0
			46	46		
4	D	24	Total	O	0	0
			24	24		

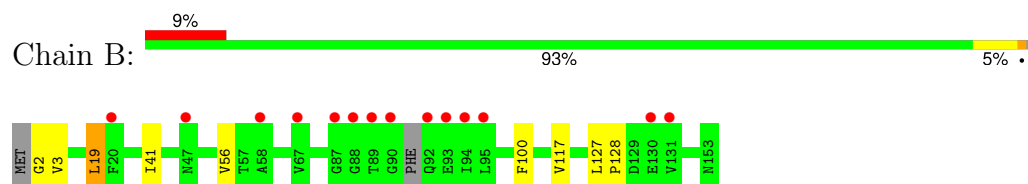
### 3 Residue-property plots [i](#)

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.

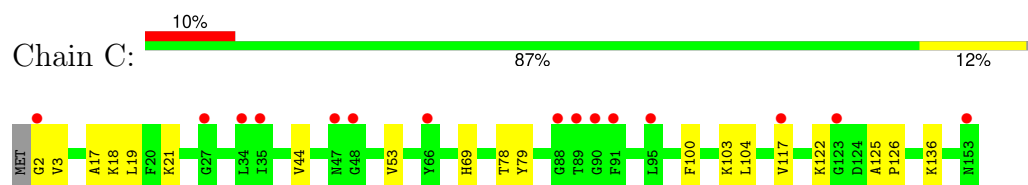
- Molecule 1: Ara h 8 allergen isoform



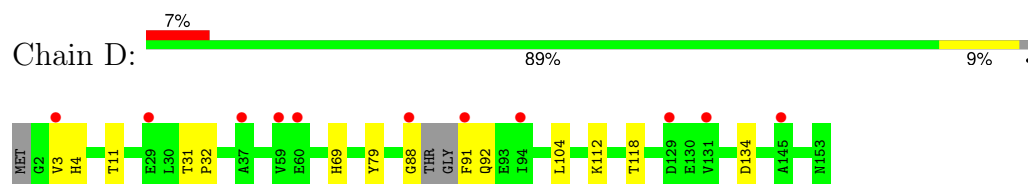
- Molecule 1: Ara h 8 allergen isoform



- Molecule 1: Ara h 8 allergen isoform



- Molecule 1: Ara h 8 allergen isoform



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	38.41Å 59.21Å 67.08Å 89.96° 89.99° 71.02°	Depositor
Resolution (Å)	31.96 – 1.95 31.96 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.1 (31.96-1.95) 99.7 (31.96-1.95)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.35 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.173 , 0.219 0.207 , 0.254	Depositor DCC
$R_{free}$ test set	2121 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	24.9	Xtriage
Anisotropy	0.218	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 36.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.54$ , $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	0.259 for h,h-k,-l 0.047 for -h,-k,l 0.056 for -h,-h+k,-l	Xtriage
Reported twinning fraction	0.365 for H, K, L 0.253 for -h,-k,l 0.204 for H, H-K, -L 0.177 for -H, -H+K, -L	Depositor
Outliers	0 of 40698 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4721	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.93% 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: SO4, NA

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	0.76	0/1154	0.83	0/1565
1	B	0.75	0/1147	0.78	0/1554
1	C	0.76	0/1154	0.80	0/1565
1	D	0.76	0/1152	0.83	0/1561
All	All	0.76	0/4607	0.81	0/6245

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	1133	0	1123	15	0
1	B	1127	0	1118	6	0
1	C	1133	0	1116	13	0
1	D	1131	0	1114	10	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	15	0	0	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	5	0	0	0	0
3	C	15	0	0	1	0
3	D	5	0	0	0	0
4	A	42	0	0	1	0
4	B	41	0	0	0	0
4	C	46	0	0	3	0
4	D	24	0	0	0	0
All	All	4721	0	4471	40	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 4.

All (40) 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:A:9:GLU:HB3	1:A:112:LYS:HZ2	1.46	0.78
1:D:88:GLY:HA3	1:D:91:PHE:HB2	1.65	0.78
1:C:19:LEU:HD23	1:C:104:LEU:HD13	1.72	0.70
1:A:9:GLU:HB3	1:A:112:LYS:NZ	2.08	0.68
1:C:17:ALA:HA	1:C:79:TYR:CE1	2.34	0.62
1:D:92:GLN:NE2	1:D:134:ASP:OD2	2.33	0.62
1:D:11:THR:HG22	1:D:112:LYS:HG2	1.83	0.60
1:B:100:PHE:CD2	1:B:117:VAL:HG22	2.37	0.60
1:A:124:ASP:O	1:A:124:ASP:OD1	2.19	0.59
1:A:125:ALA:HB2	4:A:334:HOH:O	2.01	0.59
1:D:88:GLY:HA3	1:D:91:PHE:CB	2.33	0.57
1:A:19:LEU:HD23	1:A:104:LEU:HD13	1.87	0.57
1:C:78:THR:HB	1:C:103:LYS:NZ	2.21	0.56
1:D:79:TYR:CB	1:D:104:LEU:HB2	2.37	0.55
1:A:8:GLU:HA	1:A:8:GLU:OE2	2.08	0.54
1:A:5:THR:HG22	4:C:316:HOH:O	2.08	0.53
1:A:4:HIS:HA	1:C:3:VAL:O	2.10	0.52
1:A:92:GLN:NE2	1:A:134:ASP:OD2	2.44	0.51
1:D:79:TYR:HB2	1:D:104:LEU:HB2	1.92	0.50
1:A:34:LEU:HD21	1:A:149:TYR:CD2	2.48	0.49
1:A:95:LEU:HD11	1:A:131:VAL:CG1	2.43	0.49
1:C:44:VAL:HG22	1:C:53:VAL:O	2.13	0.49
1:A:4:HIS:CE1	1:A:126:PRO:HA	2.48	0.48
1:D:3:VAL:HG13	1:D:118:THR:HG23	1.95	0.48
1:B:2:GLY:HA2	1:D:3:VAL:O	2.14	0.48
1:C:21:LYS:HD3	4:C:324:HOH:O	2.14	0.48

*Continued on next page...*



Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:VAL:O	1:D:4:HIS:HA	2.15	0.46
1:B:127:LEU:HD12	1:B:128:PRO:HD2	1.97	0.46
1:C:136:LYS:HZ3	1:C:136:LYS:HB2	1.81	0.45
1:C:100:PHE:CD2	1:C:117:VAL:HG22	2.52	0.45
1:A:3:VAL:O	1:C:2:GLY:HA2	2.18	0.44
1:A:79:TYR:HD2	1:A:104:LEU:O	2.02	0.43
1:B:19:LEU:HD12	1:B:19:LEU:O	2.19	0.43
1:C:44:VAL:CG2	1:C:53:VAL:HG12	2.48	0.43
1:C:18:LYS:HD3	3:C:203:SO4:O3	2.19	0.43
1:A:38:ILE:O	1:A:38:ILE:HG22	2.18	0.42
1:B:41:ILE:HG12	1:B:56:VAL:HG22	2.02	0.41
1:C:125:ALA:HA	1:C:126:PRO:HD2	1.94	0.41
1:C:136:LYS:HA	4:C:343:HOH:O	2.20	0.41
1:D:31:THR:N	1:D:32:PRO:CD	2.85	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	150/153 (98%)	150 (100%)	0	0	100	100
1	B	147/153 (96%)	145 (99%)	2 (1%)	0	100	100
1	C	150/153 (98%)	150 (100%)	0	0	100	100
1	D	146/153 (95%)	146 (100%)	0	0	100	100
All	All	593/612 (97%)	591 (100%)	2 (0%)	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	123/128 (96%)	123 (100%)	0	100	100
1	B	123/128 (96%)	122 (99%)	1 (1%)	79	78
1	C	123/128 (96%)	121 (98%)	2 (2%)	58	55
1	D	124/128 (97%)	123 (99%)	1 (1%)	79	78
All	All	493/512 (96%)	489 (99%)	4 (1%)	79	78

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

Mol	Chain	Res	Type
1	B	19	LEU
1	C	69	HIS
1	C	122	LYS
1	D	69	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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](#)

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 for Mogul analysis.

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	SO4	A	204	-	4,4,4	0.32	0	6,6,6	0.10	0
3	SO4	A	202	-	4,4,4	0.51	0	6,6,6	0.06	0
3	SO4	C	202	-	4,4,4	0.59	0	6,6,6	0.17	0
3	SO4	C	204	-	4,4,4	0.33	0	6,6,6	0.07	0
3	SO4	D	202	-	4,4,4	0.31	0	6,6,6	0.13	0
3	SO4	A	203	-	4,4,4	0.32	0	6,6,6	0.08	0
3	SO4	B	202	-	4,4,4	0.28	0	6,6,6	0.09	0
3	SO4	C	203	-	4,4,4	0.35	0	6,6,6	0.06	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	203	SO4	1	0

## 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	152/153 (99%)	1.07	20 (13%) 8 10	22, 31, 49, 58	0
1	B	151/153 (98%)	0.87	14 (9%) 16 19	22, 30, 45, 55	0
1	C	152/153 (99%)	0.91	15 (9%) 14 18	23, 31, 43, 56	0
1	D	150/153 (98%)	0.91	11 (7%) 22 27	22, 30, 44, 61	0
All	All	605/612 (98%)	0.94	60 (9%) 14 18	22, 31, 47, 61	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	91	PHE	7.0
1	C	91	PHE	5.7
1	A	94	ILE	4.9
1	A	89	THR	4.7
1	B	89	THR	3.9
1	D	94	ILE	3.7
1	C	47	ASN	3.7
1	A	87	GLY	3.4
1	A	88	GLY	3.3
1	B	94	ILE	3.2
1	B	47	ASN	3.1
1	B	92	GLN	3.1
1	A	58	ALA	3.1
1	D	37	ALA	3.0
1	D	91	PHE	3.0
1	C	153	ASN	3.0
1	B	67	VAL	3.0
1	B	90	GLY	2.9
1	B	95	LEU	2.9
1	C	95	LEU	2.9
1	D	131	VAL	2.9

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	C	90	GLY	2.9
1	D	60	GLU	2.8
1	C	123	GLY	2.8
1	B	130	GLU	2.7
1	A	90	GLY	2.7
1	A	33	LYS	2.6
1	A	122	LYS	2.6
1	A	47	ASN	2.5
1	D	145	ALA	2.4
1	D	59	VAL	2.4
1	C	34	LEU	2.4
1	A	125	ALA	2.4
1	D	29	GLU	2.4
1	C	35	ILE	2.3
1	C	117	VAL	2.3
1	C	89	THR	2.3
1	B	58	ALA	2.2
1	B	87	GLY	2.2
1	D	3	VAL	2.2
1	C	2	GLY	2.2
1	C	88	GLY	2.2
1	D	129	ASP	2.2
1	B	88	GLY	2.2
1	A	3	VAL	2.1
1	A	32	PRO	2.1
1	B	93	GLU	2.1
1	A	59	VAL	2.1
1	B	131	VAL	2.1
1	A	29	GLU	2.1
1	C	66	TYR	2.1
1	A	148	GLY	2.1
1	C	27	GLY	2.1
1	C	48	GLY	2.1
1	D	88	GLY	2.1
1	A	67	VAL	2.0
1	A	130	GLU	2.0
1	B	20	PHE	2.0
1	A	79	TYR	2.0
1	A	107	ALA	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](#)

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. 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	B-factors( $\text{\AA}^2$ )	Q<0.9
3	SO4	A	202	5/5	0.75	0.13	69,70,73,74	0
3	SO4	A	204	5/5	0.81	0.11	56,61,63,66	0
3	SO4	C	203	5/5	0.81	0.24	43,44,45,51	0
3	SO4	A	203	5/5	0.83	0.12	49,55,57,58	0
3	SO4	C	204	5/5	0.84	0.11	55,55,56,59	0
3	SO4	C	202	5/5	0.85	0.13	46,46,51,53	0
3	SO4	B	202	5/5	0.86	0.18	39,39,41,43	0
3	SO4	D	202	5/5	0.88	0.16	45,48,49,53	0
2	NA	B	201	1/1	0.92	0.07	37,37,37,37	0
2	NA	A	201	1/1	0.96	0.05	25,25,25,25	0
2	NA	D	201	1/1	0.97	0.05	23,23,23,23	0
2	NA	C	201	1/1	0.99	0.03	32,32,32,32	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.