



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 08:31 AM UTC

PDB ID : 5VPK / pdb_00005vpk
Title : CRYSTAL STRUCTURE OF MITE ALLERGEN DER F 1
Authors : Chruszcz, M.; Chapman, M.D.; Vailes, L.D.; Pomes, A.; Minor, W.
Deposited on : 2017-05-05
Resolution : 2.00 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
Mogul : 2022.3.0, CSD as543be (2022)
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

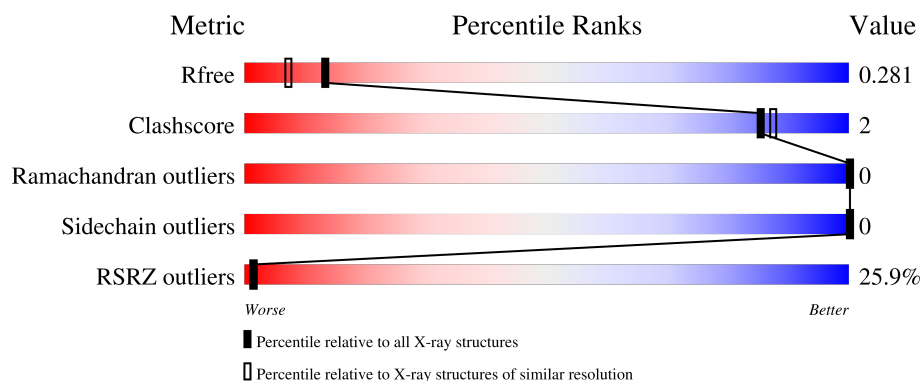
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	223	11% 93% 7%
1	B	223	12% 95% 5%
1	C	223	54% 94% 6%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5614 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 Der f 1 variant.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	223	Total	C	N	O	S	0	1	0
			1777	1108	315	342	12			
1	B	223	Total	C	N	O	S	0	1	0
			1779	1109	317	342	11			
1	C	223	Total	C	N	O	S	0	1	0
			1711	1075	293	332	11			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

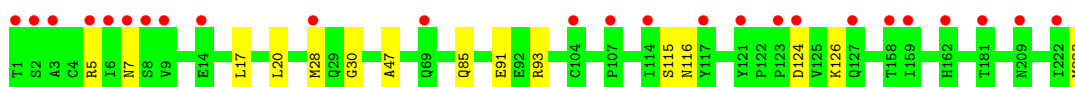
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	122	Total	O	0	0
			122	122		
4	B	118	Total	O	0	0
			118	118		
4	C	49	Total	O	0	0
			49	49		

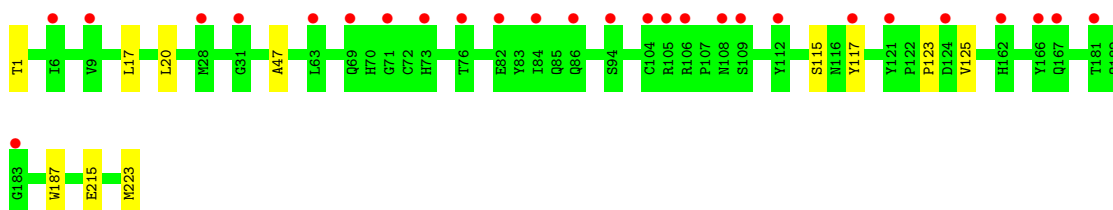
3 Residue-property plots [i](#)

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.

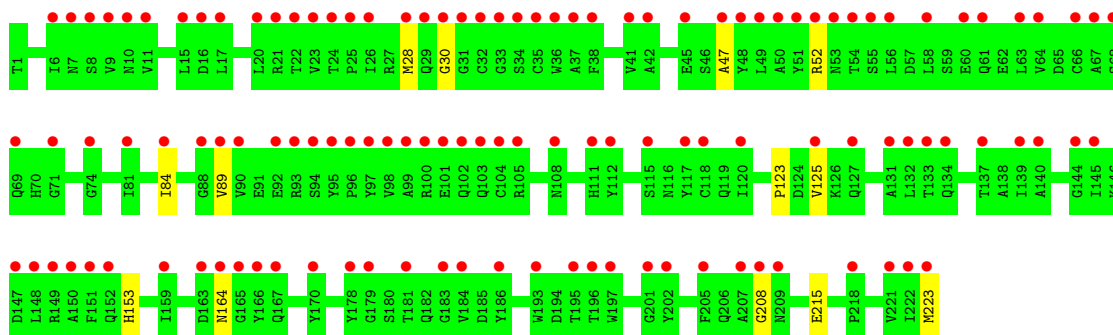
- Molecule 1: Der f 1 variant



- Molecule 1: Der f 1 variant



- Molecule 1: Der f 1 variant



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	91.16Å 91.16Å 77.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.10 – 2.00 36.10 – 2.00	Depositor EDS
% Data completeness (in resolution range)	92.4 (36.10-2.00) 92.4 (36.10-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.99 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.180 , 0.210 (Not available) , 0.281	Depositor DCC
R_{free} test set	1958 reflections (4.54%)	wwPDB-VP
Wilson B-factor (Å ²)	30.9	Xtriage
Anisotropy	0.134	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.417 for h,-k,-l	Xtriage
Reported twinning fraction	0.597 for H, K, L 0.403 for -H, K, -L	Depositor
Outliers	0 of 39868 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5614	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, SO4

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	0.80	0/1820	0.85	0/2476
1	B	0.86	0/1823	0.88	0/2481
1	C	0.73	0/1754	0.82	0/2398
All	All	0.80	0/5397	0.85	0/7355

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	1777	0	1676	9	1
1	B	1779	0	1674	6	1
1	C	1711	0	1563	9	0
2	A	25	0	0	0	0
2	B	5	0	0	0	0
3	A	14	0	13	0	0
3	B	14	0	13	0	0
4	A	122	0	0	1	0
4	B	118	0	0	0	0
4	C	49	0	0	2	0
All	All	5614	0	4939	23	1

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 (23) 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:115:SER:O	1:C:164:ASN:HB3	1.91	0.70
1:C:28:MET:HE3	1:C:30:GLY:H	1.64	0.62
1:A:91:GLU:OE2	1:A:93:ARG:NH1	2.34	0.59
1:A:5:ARG:NH2	4:A:401:HOH:O	2.38	0.57
1:C:153:HIS:NE2	4:C:301:HOH:O	2.31	0.55
1:B:125:VAL:HG23	1:B:187:TRP:CZ2	2.42	0.53
1:A:91:GLU:CD	1:A:93:ARG:NH1	2.66	0.53
1:C:28:MET:HE3	1:C:30:GLY:N	2.24	0.53
1:B:125:VAL:HG23	1:B:187:TRP:HZ2	1.75	0.50
1:C:47:ALA:HB1	1:C:223:MET:HG2	1.96	0.48
1:C:123:PRO:HB2	1:C:215:GLU:O	2.13	0.48
1:A:28[B]:MET:HE3	1:A:30:GLY:HA2	1.95	0.47
1:B:123:PRO:HB2	1:B:215:GLU:O	2.13	0.47
1:A:47:ALA:HB1	1:A:223:MET:HG2	1.96	0.47
1:B:47:ALA:HB1	1:B:223:MET:HG2	1.97	0.46
1:C:125:VAL:HG21	1:C:208:GLY:HA2	1.96	0.46
1:A:124:ASP:OD1	1:A:126:LYS:N	2.45	0.45
1:B:1:THR:HG23	1:B:117:TYR:CE2	2.51	0.45
1:C:52:ARG:NH2	4:C:307:HOH:O	2.51	0.42
1:A:85:GLN:OE1	1:A:116:ASN:HA	2.21	0.41
1:C:84:ILE:HG12	1:C:89:VAL:HG13	2.03	0.41
1:A:17:LEU:HD23	1:A:20:LEU:HD12	2.02	0.40
1:B:17:LEU:HD23	1:B:20:LEU:HD12	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:ASN:ND2	1:B:115:SER:OG[2_455]	2.06	0.14

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	222/223 (100%)	215 (97%)	7 (3%)	0	100	100
1	B	222/223 (100%)	214 (96%)	8 (4%)	0	100	100
1	C	222/223 (100%)	217 (98%)	5 (2%)	0	100	100
All	All	666/669 (100%)	646 (97%)	20 (3%)	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	191/190 (100%)	191 (100%)	0	100	100
1	B	191/190 (100%)	191 (100%)	0	100	100
1	C	176/190 (93%)	176 (100%)	0	100	100
All	All	558/570 (98%)	558 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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	85	GLN
1	A	108	ASN
1	A	153	HIS
1	A	161	GLN
1	A	209	ASN
1	B	153	HIS
1	C	206	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

8 ligands 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
2	SO4	A	302	-	4,4,4	0.42	0	6,6,6	0.14	0
2	SO4	A	304	-	4,4,4	0.46	0	6,6,6	0.10	0
3	NAG	B	301	1	14,14,15	0.58	0	17,19,21	1.55	5 (29%)
2	SO4	B	302	-	4,4,4	0.41	0	6,6,6	0.21	0
3	NAG	A	303	1	14,14,15	0.34	0	17,19,21	1.70	2 (11%)
2	SO4	A	301	-	4,4,4	0.45	0	6,6,6	0.11	0
2	SO4	A	305	-	4,4,4	0.48	0	6,6,6	0.29	0
2	SO4	A	306	-	4,4,4	0.43	0	6,6,6	0.11	0

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	NAG	B	301	1	-	0/6/23/26	0/1/1/1
3	NAG	A	303	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	303	NAG	C1-O5-C5	5.01	118.91	112.19
3	A	303	NAG	C1-C2-N2	-3.17	105.43	110.43
3	B	301	NAG	C1-O5-C5	2.75	115.88	112.19
3	B	301	NAG	C1-C2-N2	-2.55	106.42	110.43
3	B	301	NAG	C4-C3-C2	-2.49	107.36	111.02
3	B	301	NAG	C6-C5-C4	-2.27	107.45	113.02
3	B	301	NAG	C3-C4-C5	2.08	114.00	110.23

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	303	NAG	C4-C5-C6-O6
3	A	303	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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, 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	223/223 (100%)	0.99	25 (11%)	10 9	17, 35, 50, 62	1 (0%)
1	B	223/223 (100%)	0.87	27 (12%)	8 8	18, 34, 54, 77	1 (0%)
1	C	223/223 (100%)	2.21	121 (54%)	0 0	29, 52, 88, 107	1 (0%)
All	All	669/669 (100%)	1.36	173 (25%)	1 1	17, 39, 74, 107	3 (0%)

All (173) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	9	VAL	6.1
1	C	223	MET	4.8
1	C	108	ASN	4.7
1	C	53	ASN	4.7
1	C	56	LEU	4.7
1	C	37	ALA	4.4
1	C	54	THR	4.2
1	C	21[A]	ARG	4.1
1	C	98	VAL	4.1
1	C	104	CYS	4.0
1	C	7	ASN	4.0
1	C	51	TYR	4.0
1	C	99	ALA	3.9
1	B	108	ASN	3.9
1	C	101	GLU	3.9
1	C	49	LEU	3.9
1	C	67	ALA	3.9
1	C	97	TYR	3.8
1	C	164	ASN	3.8
1	C	89	VAL	3.7
1	A	209	ASN	3.6
1	C	31	GLY	3.6
1	A	5	ARG	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	31	GLY	3.6
1	B	183	GLY	3.6
1	C	202	TYR	3.6
1	C	74	GLY	3.5
1	B	166	TYR	3.4
1	C	103	GLN	3.4
1	C	30	GLY	3.4
1	C	96	PRO	3.4
1	C	93	ARG	3.4
1	A	8	SER	3.4
1	C	42	ALA	3.4
1	C	140	ALA	3.4
1	C	148	LEU	3.3
1	C	193	TRP	3.3
1	C	47	ALA	3.3
1	B	167	GLN	3.3
1	B	104	CYS	3.2
1	C	25	PRO	3.2
1	C	48	TYR	3.2
1	C	63	LEU	3.2
1	A	2	SER	3.2
1	C	6	ILE	3.2
1	C	58	LEU	3.2
1	C	222	ILE	3.2
1	C	36	TRP	3.1
1	C	90	VAL	3.1
1	C	131	ALA	3.1
1	A	181	THR	3.1
1	C	55	SER	3.1
1	C	60	GLU	3.0
1	C	28	MET	3.0
1	C	8	SER	3.0
1	C	139	ILE	3.0
1	C	95	TYR	3.0
1	C	100	ARG	3.0
1	B	6	ILE	3.0
1	A	9	VAL	3.0
1	C	10	ASN	2.9
1	C	11	VAL	2.9
1	A	28[A]	MET	2.9
1	C	32	CYS	2.9
1	C	52	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	105	ARG	2.9
1	A	124	ASP	2.8
1	C	195	THR	2.8
1	C	71	GLY	2.8
1	C	17	LEU	2.8
1	B	181	THR	2.8
1	C	16	ASP	2.8
1	C	33	GLY	2.8
1	C	166	TYR	2.8
1	C	178	TYR	2.8
1	C	24	THR	2.8
1	C	149	ARG	2.8
1	C	133	THR	2.8
1	C	26	ILE	2.8
1	C	197	TRP	2.8
1	B	112	TYR	2.7
1	C	137	THR	2.7
1	A	107	PRO	2.7
1	C	88	GLY	2.7
1	B	86	GLN	2.7
1	C	41	VAL	2.7
1	C	94	SER	2.7
1	C	145	ILE	2.6
1	B	71	GLY	2.6
1	C	15	LEU	2.6
1	C	186	TYR	2.6
1	C	64	VAL	2.5
1	C	84	ILE	2.5
1	C	165	GLY	2.5
1	C	147	ASP	2.5
1	A	117	TYR	2.5
1	A	121	TYR	2.5
1	C	117	TYR	2.5
1	C	184	VAL	2.5
1	A	69	GLN	2.5
1	C	102	GLN	2.5
1	B	109	SER	2.5
1	C	50	ALA	2.5
1	B	82	GLU	2.4
1	C	127	GLN	2.4
1	C	118	CYS	2.4
1	C	205	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	6	ILE	2.4
1	C	221	VAL	2.4
1	C	20	LEU	2.4
1	C	183	GLY	2.4
1	C	134	GLN	2.4
1	C	23	VAL	2.4
1	C	120	ILE	2.4
1	C	34	SER	2.4
1	C	66	CYS	2.4
1	C	22	THR	2.4
1	C	111	HIS	2.4
1	A	3	ALA	2.4
1	C	152	GLN	2.4
1	C	45	GLU	2.3
1	C	207	ALA	2.3
1	C	68	SER	2.3
1	C	159	ILE	2.3
1	C	69	GLN	2.3
1	C	209	ASN	2.3
1	C	201	GLY	2.3
1	B	94	SER	2.3
1	B	162[A]	HIS	2.3
1	A	158	THR	2.3
1	B	9	VAL	2.2
1	A	7	ASN	2.2
1	C	35	CYS	2.2
1	C	92	GLU	2.2
1	C	170	TYR	2.2
1	C	151	PHE	2.2
1	C	181	THR	2.2
1	A	127	GLN	2.2
1	C	29	GLN	2.2
1	C	112	TYR	2.2
1	C	218	PRO	2.2
1	B	117	TYR	2.2
1	B	28	MET	2.2
1	A	114	ILE	2.1
1	C	61	GLN	2.1
1	A	162	HIS	2.1
1	C	132	LEU	2.1
1	A	123	PRO	2.1
1	B	76	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	167	GLN	2.1
1	C	125	VAL	2.1
1	C	179	GLY	2.1
1	B	63	LEU	2.1
1	A	104	CYS	2.1
1	C	115	SER	2.1
1	A	1	THR	2.1
1	A	159	ILE	2.1
1	C	144	GLY	2.1
1	C	38	PHE	2.1
1	C	150	ALA	2.1
1	A	14	GLU	2.1
1	B	69	GLN	2.1
1	C	105	ARG	2.1
1	C	208	GLY	2.1
1	A	222	ILE	2.1
1	C	81	ILE	2.1
1	B	124	ASP	2.0
1	B	73	HIS	2.0
1	C	196	THR	2.0
1	B	84	ILE	2.0
1	C	163	ASP	2.0
1	B	106	ARG	2.0
1	B	121	TYR	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](#)

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, 95th 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
2	SO4	A	302	5/5	0.58	0.14	103,106,108,109	0
2	SO4	A	301	5/5	0.68	0.13	84,86,88,89	0
3	NAG	B	301	14/15	0.72	0.15	43,50,56,56	0
2	SO4	B	302	5/5	0.73	0.18	75,76,83,84	0
3	NAG	A	303	14/15	0.73	0.13	46,55,60,63	0
2	SO4	A	305	5/5	0.73	0.16	75,75,77,77	0
2	SO4	A	304	5/5	0.79	0.12	82,82,83,87	0
2	SO4	A	306	5/5	0.83	0.15	77,77,80,86	0

6.5 Other polymers [i](#)

There are no such residues in this entry.