



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:27 PM UTC

PDB ID : 4HGM / pdb_00004hgm
Title : Shark IgNAR Variable Domain
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Deposited on : 2012-10-08
Resolution : 2.34 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

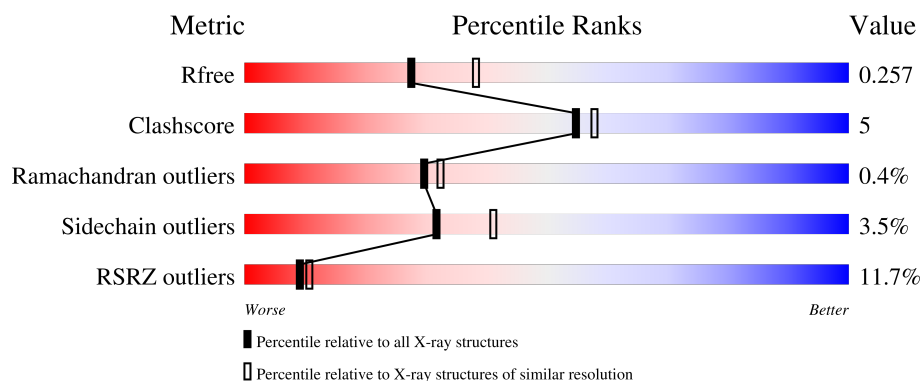
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.34 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	3031 (2.36-2.32)
Clashscore	190562	3127 (2.36-2.32)
Ramachandran outliers	187476	3095 (2.36-2.32)
Sidechain outliers	187428	3095 (2.36-2.32)
RSRZ outliers	180081	3033 (2.36-2.32)

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	112	4% 79% 11% 8%
2	B	585	11% 70% 10% 18%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4577 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.

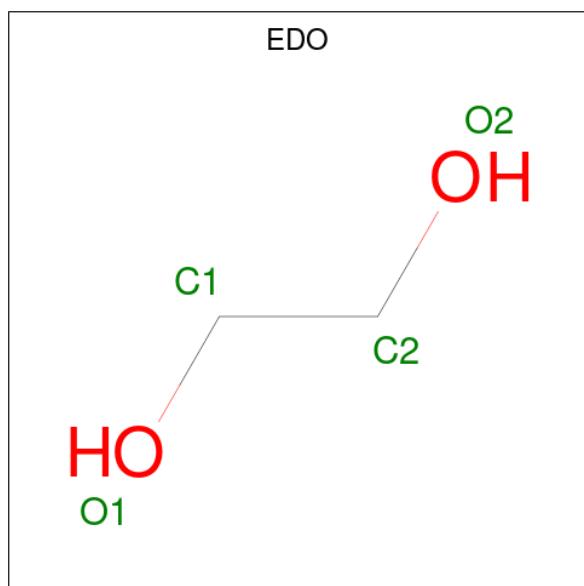
- Molecule 1 is a protein called Shark V-NAR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	103	Total	C	N	O	S	0	0	0
			791	494	130	164	3			

- Molecule 2 is a protein called Serum albumin.

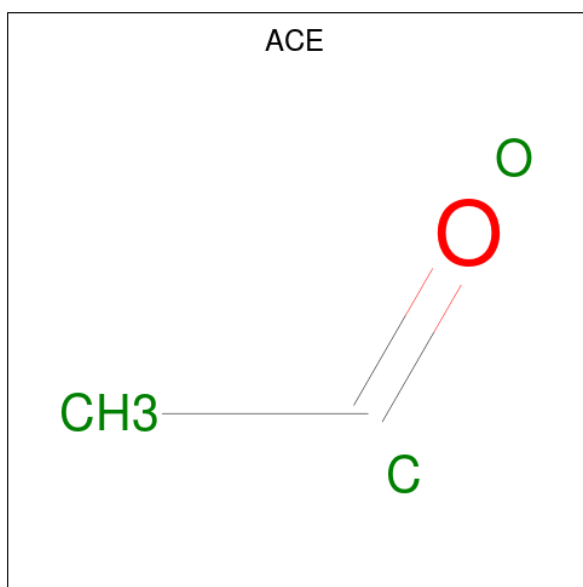
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	478	Total	C	N	O	S	0	0	0
			3707	2342	632	697	36			

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is ACETYL GROUP (CCD ID: ACE) (formula: C₂H₄O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			3	2	1		

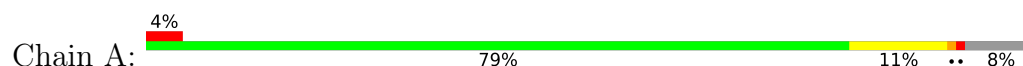
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	22	Total	O	0	0
			22	22		
5	B	50	Total	O	0	0
			50	50		

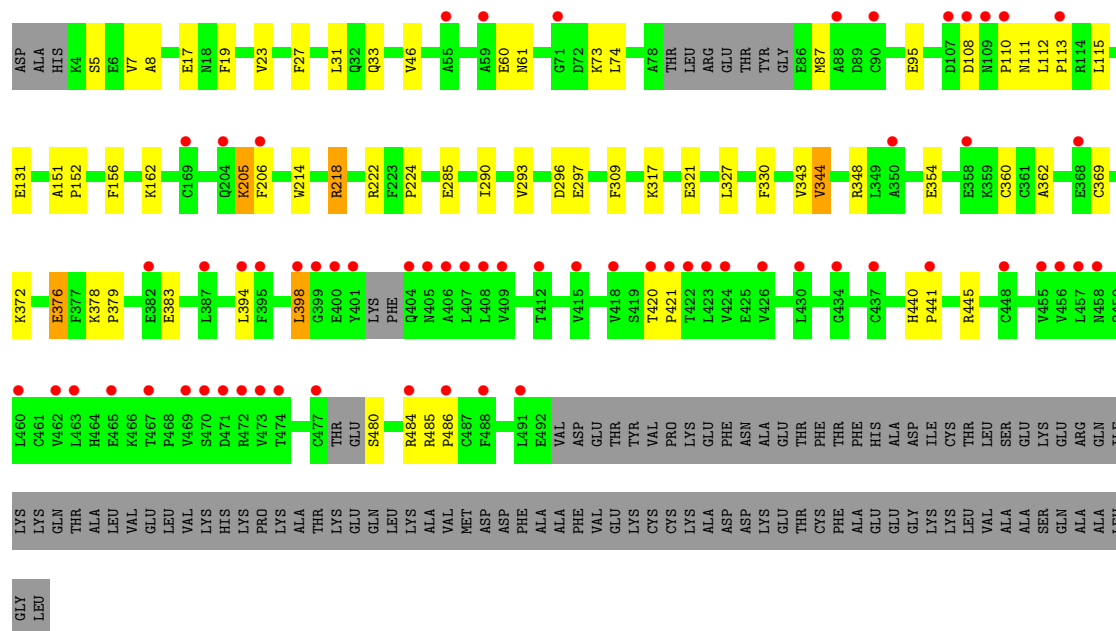
3 Residue-property plots [i](#)

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: Shark V-NAR



• Molecule 2: Serum albumin



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	131.34Å 131.34Å 74.52Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.91 – 2.34 37.91 – 2.34	Depositor EDS
% Data completeness (in resolution range)	95.4 (37.91-2.34) 99.5 (37.91-2.34)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 2.34Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, R_{free}	0.221 , 0.259 0.221 , 0.257	Depositor DCC
R_{free} test set	1604 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	48.4	Xtriage
Anisotropy	0.198	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 51.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.038 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4577	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% 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: ACE, EDO

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	0.36	0/807	0.69	0/1096
2	B	0.34	0/3779	0.72	0/5111
All	All	0.34	0/4586	0.72	0/6207

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	791	0	766	8	0
2	B	3707	0	3546	34	0
3	B	4	0	6	0	0
4	B	3	0	3	0	0
5	A	22	0	0	0	0
5	B	50	0	0	0	0
All	All	4577	0	4321	42	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 (42) 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:485:ARG:HB3	2:B:486:PRO:HD3	1.70	0.73
2:B:222:ARG:C	2:B:224:PRO:HD3	2.28	0.58
2:B:317:LYS:HE2	2:B:321:GLU:OE2	2.04	0.57
2:B:378:LYS:HB2	2:B:379:PRO:HD3	1.86	0.56
1:A:89:ASN:O	1:A:90:ILE:HG22	2.06	0.56
1:A:76:GLU:H	1:A:76:GLU:CD	2.14	0.56
2:B:112:LEU:HB3	2:B:113:PRO:HD2	1.90	0.53
2:B:156:PHE:CE1	2:B:285:GLU:HB3	2.45	0.52
2:B:344:VAL:O	2:B:348:ARG:HG3	2.10	0.51
2:B:224:PRO:HD2	2:B:296:ASP:HB3	1.93	0.51
2:B:327:LEU:HD11	2:B:354:GLU:HG3	1.94	0.49
2:B:376:GLU:O	2:B:379:PRO:HD2	2.12	0.49
2:B:480:SER:O	2:B:484:ARG:HG3	2.13	0.49
2:B:441:PRO:O	2:B:445:ARG:HG3	2.13	0.49
1:A:11:SER:HA	1:A:101:GLU:O	2.12	0.49
2:B:5:SER:HB3	2:B:8:ALA:HB3	1.95	0.48
2:B:218:ARG:HD3	2:B:218:ARG:C	2.38	0.48
2:B:420:THR:N	2:B:421:PRO:HD2	2.28	0.48
2:B:372:LYS:HE2	2:B:372:LYS:HA	1.96	0.47
2:B:394:LEU:O	2:B:398:LEU:HB2	2.14	0.47
2:B:218:ARG:HD3	2:B:218:ARG:O	2.14	0.47
2:B:214:TRP:CD1	2:B:343:VAL:HG11	2.49	0.46
1:A:29:TYR:CD1	1:A:89:ASN:HB3	2.51	0.46
2:B:60:GLU:O	2:B:61:ASN:HB2	2.17	0.45
2:B:31:LEU:HD22	2:B:74:LEU:HD22	1.99	0.45
1:A:36:TRP:CZ3	1:A:83:CYS:HB3	2.52	0.45
2:B:156:PHE:HE1	2:B:285:GLU:HB3	1.81	0.45
2:B:360:CYS:C	2:B:362:ALA:H	2.25	0.44
2:B:290:ILE:O	2:B:293:VAL:HG12	2.18	0.44
1:A:42:GLY:O	1:A:43:SER:HB2	2.18	0.44
2:B:46:VAL:HG23	2:B:73:LYS:HG2	2.00	0.43
2:B:440:HIS:HB3	2:B:441:PRO:HD2	2.00	0.43
1:A:1:THR:HG23	1:A:27:THR:HB	2.00	0.43
2:B:309:PHE:CZ	2:B:330:PHE:HA	2.54	0.43
1:A:39:GLN:HG2	1:A:43:SER:HA	2.00	0.42
2:B:115:LEU:HD23	2:B:115:LEU:HA	1.93	0.42
2:B:27:PHE:CD1	2:B:74:LEU:HD21	2.55	0.41
2:B:151:ALA:N	2:B:152:PRO:HD2	2.36	0.41
2:B:19:PHE:O	2:B:23:VAL:HG23	2.21	0.41
2:B:360:CYS:HB3	2:B:369:CYS:HB3	1.93	0.40
2:B:205:LYS:HB3	2:B:206:PHE:HD2	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:131:GLU:OE2	2:B:162:LYS:HE3	2.21	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	101/112 (90%)	97 (96%)	3 (3%)	1 (1%)	12	11
2	B	470/585 (80%)	447 (95%)	22 (5%)	1 (0%)	43	50
All	All	571/697 (82%)	544 (95%)	25 (4%)	2 (0%)	30	32

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	90	ILE
2	B	110	PRO

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	90/97 (93%)	87 (97%)	3 (3%)	33	43
2	B	392/511 (77%)	378 (96%)	14 (4%)	31	40
All	All	482/608 (79%)	465 (96%)	17 (4%)	32	41

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	VAL
1	A	76	GLU
1	A	90	ILE
2	B	7	VAL
2	B	17	GLU
2	B	33	GLN
2	B	87	MET
2	B	95	GLU
2	B	108	ASP
2	B	111	ASN
2	B	205	LYS
2	B	218	ARG
2	B	297	GLU
2	B	344	VAL
2	B	376	GLU
2	B	383	GLU
2	B	398	LEU

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

Mol	Chain	Res	Type
2	B	29	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](#)

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](#)

2 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	$\# Z > 2$
4	ACE	B	602	-	1,2,2	0.89	0	0,1,1	-	-
3	EDO	B	601	-	3,3,3	0.33	0	2,2,2	0.48	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	EDO	B	601	-	-	0/1/1/1	-

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.

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	103/112 (91%)	0.22	4 (3%) 43 49	32, 47, 65, 80	0
2	B	478/585 (81%)	0.80	64 (13%) 7 8	33, 58, 155, 173	0
All	All	581/697 (83%)	0.70	68 (11%) 9 11	32, 55, 152, 173	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	460	LEU	6.2
2	B	109	ASN	6.0
2	B	477	CYS	4.1
2	B	110	PRO	4.1
2	B	462	VAL	4.0
2	B	457	LEU	3.9
2	B	407	LEU	3.6
2	B	488	PHE	3.5
2	B	458	ASN	3.5
1	A	44	SER	3.4
2	B	471	ASP	3.4
2	B	409	VAL	3.4
2	B	415	VAL	3.4
2	B	401	TYR	3.3
2	B	456	VAL	3.3
2	B	113	PRO	3.2
2	B	423	LEU	3.2
2	B	59	ALA	3.2
2	B	441	PRO	3.2
2	B	398	LEU	3.1
2	B	467	THR	3.1
2	B	169	CYS	3.1
2	B	463	LEU	3.1
2	B	474	THR	3.1

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Mol	Chain	Res	Type	RSRZ
2	B	405	ASN	3.1
2	B	90	CYS	3.0
2	B	406	ALA	3.0
2	B	491	LEU	3.0
2	B	394	LEU	2.9
2	B	404	GLN	2.9
2	B	399	GLY	2.8
1	A	45	ASN	2.8
2	B	88	ALA	2.7
2	B	420	THR	2.7
2	B	472	ARG	2.6
2	B	107	ASP	2.6
2	B	473	VAL	2.5
2	B	465	GLU	2.5
2	B	426	VAL	2.5
2	B	455	VAL	2.4
2	B	430	LEU	2.4
2	B	470	SER	2.4
2	B	408	LEU	2.4
2	B	108	ASP	2.4
2	B	484	ARG	2.3
2	B	387	LEU	2.3
2	B	55	ALA	2.3
2	B	71	GLY	2.3
2	B	424	VAL	2.3
2	B	395	PHE	2.2
2	B	418	VAL	2.2
2	B	469	VAL	2.2
2	B	206	PHE	2.2
2	B	448	CYS	2.2
2	B	412	THR	2.2
2	B	350	ALA	2.1
2	B	358	GLU	2.1
2	B	382	GLU	2.1
1	A	71	SER	2.1
2	B	422	THR	2.1
2	B	437	CYS	2.1
2	B	368	GLU	2.1
2	B	486	PRO	2.1
2	B	434	GLY	2.1
2	B	400	GLU	2.0
2	B	204	GLN	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	10	LEU	2.0
2	B	421	PRO	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
3	EDO	B	601	4/4	0.82	0.17	55,56,59,64	0
4	ACE	B	602	3/3	0.82	0.14	46,46,48,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.