

Full wwPDB X-ray Structure Validation Report (i)

Jul 24, 2023 – 04:23 pm BST

PDB ID : 8AON

Title : Oxidoreductase fragment of human QSOX1 in complex with a Fab fragment

of a humanized anti-QSOX1 antibody

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Deposited on : 2022-08-08

Resolution : 2.10 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 (i) 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 (1)) were used in the production of this report:

 $Mol Probity \quad : \quad 4.02b\text{-}467$

Mogul : 1.8.4, CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.34

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

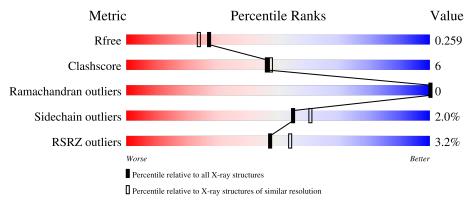
Validation Pipeline (wwPDB-VP) : 2.34

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$

The reported resolution of this entry is 2.10 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	Similar resolution
Metric	$(\# ext{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
R_{free}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 <=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	B00A	244	94%	
2	В	225	86%	10% •
3	С	214	6% 85%	15%



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5278 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 Oxidoreductase fragment of human QSOX1.

\mathbf{Mol}	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	B00A	234	Total 1848	C 1185	N 326	O 330	S 7	0	6	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B00A	29	GLY	-	expression tag	UNP A8K4C2
B00A	30	SER	-	expression tag	UNP A8K4C2
B00A	31	HIS	-	expression tag	UNP A8K4C2
B00A	32	MET	-	expression tag	UNP A8K4C2

• Molecule 2 is a protein called Heavy chain.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	В	216	Total 1596	C 1011	N 262	O 316	S 7	0	1	0

• Molecule 3 is a protein called Light chain.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
3	С	214	Total 1617	C 1007	N 271	O 333	S 6	0	1	0

• Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).





Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
4	B00A	1	Total 6	C 3	O 3	0	0

• Molecule 5 is water.

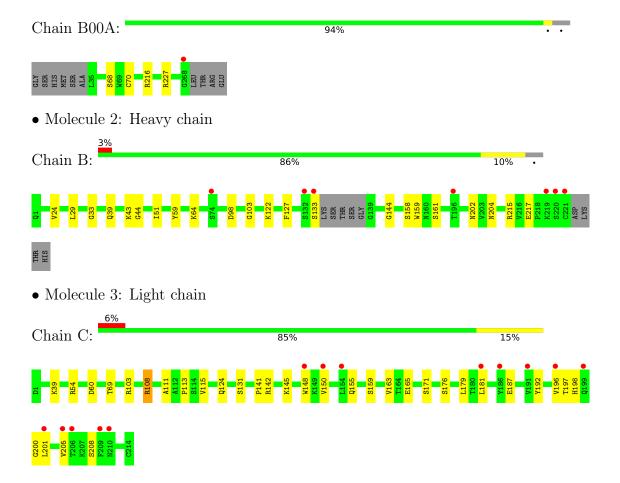
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B00A	87	Total O 87 87	0	0
5	В	65	Total O 65 65	0	0
5	С	59	Total O 59 59	0	0



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: Oxidoreductase fragment of human QSOX1





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	56.17Å 98.72Å 63.09Å	Donositor
a, b, c, α , β , γ	90.00° 101.19° 90.00°	Depositor
Resolution (Å)	49.36 - 2.10	Depositor
rtesolution (A)	49.36 - 2.10	EDS
% Data completeness	99.6 (49.36-2.10)	Depositor
(in resolution range)	99.6 (49.36-2.10)	EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.23 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.20_4459	Depositor
D D.	0.198 , 0.259	Depositor
R, R_{free}	0.197 , 0.259	DCC
R_{free} test set	3930 reflections (10.01%)	wwPDB-VP
Wilson B-factor (Å ²)	46.0	Xtriage
Anisotropy	0.319	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.33, 50.7	EDS
L-test for twinning ²	$ < L >=0.50, < L^2>=0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5278	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.16% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

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: GOL

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		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	B00A	0.42	0/1914	0.63	0/2606	
2	В	0.42	0/1638	0.61	0/2233	
3	С	0.40	0/1655	0.61	0/2258	
All	All	0.41	0/5207	0.62	0/7097	

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	B00A	1848	0	0	0	0
2	В	1596	0	1538	15	0
3	С	1617	0	1518	31	0
4	B00A	6	0	0	0	0
5	В	65	0	0	3	1
5	B00A	87	0	0	0	1
5	С	59	0	0	6	0
All	All	5278	0	3056	45	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 6.



All (45) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
2:B:161:SER:H	2:B:202:ASN:HD21	1.29	0.81
3:C:115:VAL:HG21	3:C:196:VAL:HG21	1.76	0.67
2:B:122:LYS:NZ	5:B:304:HOH:O	2.28	0.66
2:B:98[A]:ASP:OD2	5:B:301:HOH:O	2.12	0.66
3:C:103:ARG:NH2	3:C:165:GLU:OE2	2.30	0.64
3:C:201:LEU:HD12	3:C:205:VAL:HG13	1.83	0.60
3:C:142:ARG:NH2	5:C:304:HOH:O	2.34	0.60
3:C:124:GLN:HE22	3:C:131:SER:H	1.49	0.59
2:B:127:PHE:CE2	3:C:124:GLN:HG3	2.41	0.55
2:B:43:LYS:HA	2:B:43:LYS:HE2	1.88	0.55
3:C:108:ARG:HD2	3:C:171:SER:HB2	1.89	0.53
2:B:215:ARG:NH1	2:B:217:GLU:OE1	2.42	0.52
3:C:39:LYS:NZ	5:C:308:HOH:O	2.44	0.51
2:B:39:GLN:HG3	2:B:44:GLY:O	2.10	0.50
3:C:69:THR:HG22	5:C:340:HOH:O	2.11	0.50
2:B:161:SER:H	2:B:202:ASN:ND2	2.04	0.49
3:C:150:VAL:HG22	3:C:155:GLN:HE21	1.78	0.49
3:C:142:ARG:NH1	5:C:306:HOH:O	2.43	0.48
3:C:145:LYS:O	3:C:196:VAL:HA	2.13	0.48
2:B:161:SER:OG	5:B:302:HOH:O	2.20	0.48
3:C:54:ARG:HH11	3:C:54:ARG:HG3	1.78	0.47
2:B:59:TYR:HB2	2:B:64:LYS:HG2	1.95	0.47
3:C:103:ARG:HH11	3:C:103:ARG:HB2	1.80	0.47
3:C:150:VAL:HG13	3:C:155:GLN:HG3	1.95	0.47
3:C:113:PRO:HD3	3:C:198:HIS:CD2	2.52	0.45
3:C:54:ARG:HG3	3:C:54:ARG:NH1	2.33	0.45
3:C:113:PRO:HD2	3:C:201:LEU:HD23	1.99	0.44
3:C:198:HIS:CD2	3:C:200:GLY:H	2.35	0.44
3:C:148:TRP:CD2	3:C:179:LEU:HD12	2.52	0.44
3:C:179:LEU:HD22	3:C:181:LEU:HG	1.99	0.44
2:B:158:SER:HB2	2:B:202:ASN:HB2	2.00	0.43
3:C:54:ARG:NE	3:C:60:ASP:HA	2.33	0.43
3:C:163:VAL:HG21	5:C:306:HOH:O	2.18	0.43
3:C:108:ARG:NH1	3:C:111:ALA:HB2	2.33	0.43
3:C:141:PRO:O	3:C:198:HIS:HE1	2.00	0.43
2:B:144:GLY:HA2	2:B:159:TRP:CH2	2.54	0.43
3:C:103:ARG:HG2	5:C:301:HOH:O	2.19	0.42
3:C:113:PRO:HD2	3:C:201:LEU:CD2	2.49	0.42
3:C:108:ARG:HH12	3:C:111:ALA:HB2	1.85	0.42
2:B:33:GLY:HA3	2:B:51:ILE:O	2.20	0.42

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Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
2:B:24:VAL:HG21	2:B:29:LEU:HD21	2.02	0.41
3:C:148:TRP:HB3	3:C:179:LEU:HD12	2.01	0.41
2:B:98[A]:ASP:CG	2:B:103:GLY:H	2.23	0.41
3:C:192:TYR:O	3:C:208:SER:HA	2.21	0.41
3:C:148:TRP:CE3	3:C:179:LEU:HD12	2.56	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	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
5:B00A:401:HOH:O	5:B:321:HOH:O[2_345]	2.14	0.06

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	Perce	$_{ m ntiles}$
1	B00A	238/244 (98%)	234 (98%)	4 (2%)	0	100	100
2	В	$213/225\ (95\%)$	209 (98%)	4 (2%)	0	100	100
3	С	213/214 (100%)	202 (95%)	11 (5%)	0	100	100
All	All	664/683 (97%)	645 (97%)	19 (3%)	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	B00A	191/199 (96%)	187 (98%)	4 (2%)	53 59
2	В	176/187 (94%)	174 (99%)	2 (1%)	73 79
3	С	179/187 (96%)	174 (97%)	5 (3%)	43 47
All	All	546/573 (95%)	535 (98%)	11 (2%)	55 60

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

Mol	Chain	Res	Type
1	B00A	68	SER
1	B00A	70	CYS
1	B00A	216	ARG
1	B00A	227	ARG
2	В	133	SER
2	В	204	ASN
3	С	108	ARG
3	С	159	SER
3	С	176	SER
3	С	187	GLU
3	С	197	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
2	В	81	GLN
2	В	202	ASN
2	В	204	ASN
3	С	79	GLN
3	С	124	GLN
3	С	198	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand is 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	B	ond leng	$_{ m gths}$	В	ond ang	gles
IVIOI	туре	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
4	GOL	B00A	301	-	5,5,5	0.80	0	5,5,5	1.17	1 (20%)

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
4	GOL	B00A	301	-	-	4/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
4	B00A	301	GOL	C3-C2-C1	-2.26	102.94	111.70

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B00A	301	GOL	O1-C1-C2-O2
4	B00A	301	GOL	O1-C1-C2-C3
4	B00A	301	GOL	C1-C2-C3-O3
4	B00A	301	GOL	O2-C2-C3-O3



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 (i)

6.1 Protein, DNA and RNA chains (i)

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^{th} 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 $>$	#RSR2	Z>2	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q<0.9
1	B00A	234/244~(95%)	-0.02	1 (0%) 92	93	37, 51, 70, 87	0
2	В	216/225 (96%)	0.13	7 (3%) 47	54	37, 52, 75, 125	0
3	С	214/214 (100%)	0.32	13 (6%) 2	1 26	32, 53, 97, 120	0
All	All	664/683 (97%)	0.14	21 (3%) 4	7 54	32, 52, 86, 125	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	С	154	LEU	5.2
2	В	221	CYS	5.0
2	В	196	THR	4.9
1	B00A	268	GLY	3.6
2	В	132	SER	3.3
2	В	220	SER	3.3
3	С	206	THR	3.1
3	С	205	VAL	2.9
3	С	148	TRP	2.9
3	С	181	LEU	2.8
3	С	191	VAL	2.6
2	В	219	LYS	2.6
3	С	150	VAL	2.5
3	С	210	ASN	2.5
3	С	201	LEU	2.4
3	С	199	GLN	2.3
3	С	186	TYR	2.2
2	В	74	SER	2.2
3	С	196	VAL	2.1
2	В	133	SER	2.1
3	С	209	PHE	2.1



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^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
4	GOL	B00A	301	6/6	0.82	0.24	58,63,66,66	0

6.5 Other polymers (i)

There are no such residues in this entry.

