



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 8, 2024 – 10:28 AM EST

PDB ID : 2HRA  
Title : Crystal structures of the interacting domains from yeast glutamyl-tRNA synthetase and tRNA aminoacylation and nuclear export cofactor Arc1p reveal a novel function for an old fold  
Authors : Simader, H.; Hothorn, M.; Suck, D.  
Deposited on : 2006-07-20  
Resolution : 1.90 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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 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) : 2.36

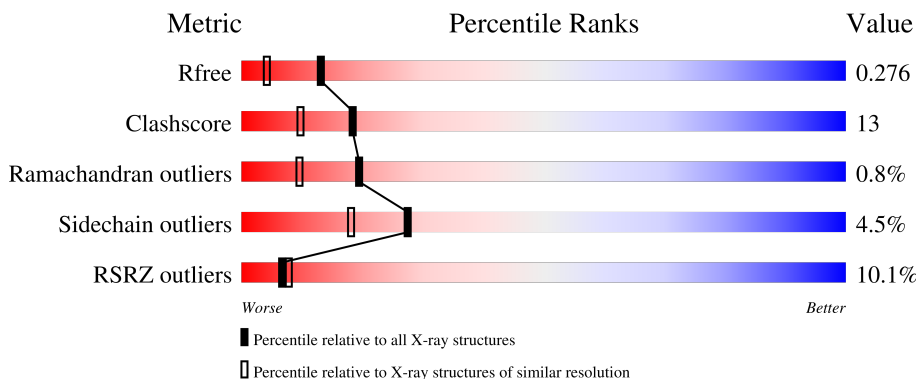
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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))
$R_{free}$	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	209	
1	B	209	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	IOD	A	402	-	-	X	-
2	IOD	A	424	-	-	X	-
2	IOD	B	414	-	-	X	-
2	IOD	B	419	-	-	X	-
2	IOD	B	421	-	-	X	-
2	IOD	B	426	-	-	X	-
2	IOD	B	440	-	-	X	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3044 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 Glutamyl-tRNA synthetase, cytoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	180	1376	881	231	259	5	0	0	0
1	B	178	1353	867	227	255	4	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	cloning artifact	UNP P46655
A	2	HIS	-	cloning artifact	UNP P46655
B	1	GLY	-	cloning artifact	UNP P46655
B	2	HIS	-	cloning artifact	UNP P46655

- Molecule 2 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	20	Total 20	I 20	0	0
2	B	24	Total 24	I 24	0	0

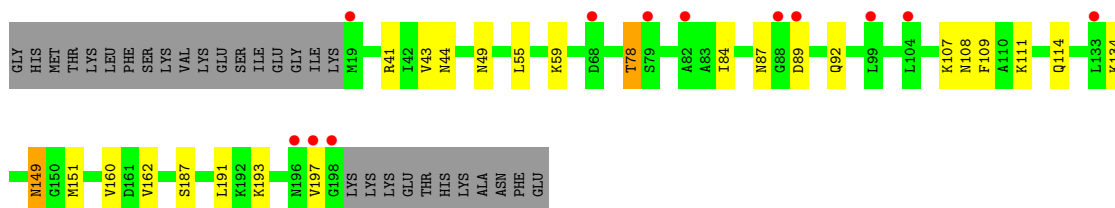
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	147	Total 147	O 147	0	0
3	B	124	Total 124	O 124	0	0

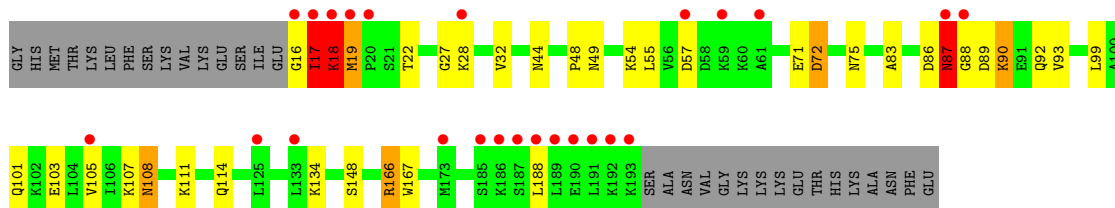
### 3 Residue-property plots [i](#)

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: Glutamyl-tRNA synthetase, cytoplasmic



- Molecule 1: Glutamyl-tRNA synthetase, cytoplasmic



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	51.96Å 107.11Å 167.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	80.00 – 1.90 45.04 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (80.00-1.90) 99.9 (45.04-1.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.66 (at 1.89Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.214 , 0.262 0.228 , 0.276	Depositor DCC
$R_{free}$ test set	1892 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	26.5	Xtrriage
Anisotropy	0.085	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 57.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3044	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.32% 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: IOD

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	0.63	0/1397	0.69	1/1889 (0.1%)
1	B	0.66	0/1374	0.72	0/1860
All	All	0.64	0/2771	0.71	1/3749 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	41	ARG	NE-CZ-NH1	5.15	122.87	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	17	ILE	Peptide
1	B	18	LYS	Peptide
1	B	87	ASN	Peptide

## 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	1376	0	1422	27	0
1	B	1353	0	1383	41	0
2	A	20	0	0	11	0
2	B	24	0	0	19	0
3	A	147	0	0	7	0
3	B	124	0	0	14	0
All	All	3044	0	2805	74	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 13.

All (74) 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:92:GLN:NE2	3:A:588:HOH:O	1.70	1.22
2:B:414:IOD:I	3:B:478:HOH:O	2.40	1.07
1:B:83:ALA:HB3	2:B:440:IOD:I	2.28	1.04
1:B:103:GLU:OE2	2:B:432:IOD:I	2.45	1.03
1:B:86:ASP:O	1:B:87:ASN:HB2	1.55	1.03
1:B:105:VAL:HG12	3:B:565:HOH:O	1.61	1.01
2:B:407:IOD:I	2:B:421:IOD:I	3.19	1.00
2:B:426:IOD:I	3:B:524:HOH:O	2.62	0.87
1:B:111:LYS:HG2	2:B:419:IOD:I	2.45	0.86
1:B:90:LYS:HE2	2:B:421:IOD:I	2.47	0.84
1:B:17:ILE:O	1:B:19:MET:CB	2.25	0.84
1:A:78:THR:HG21	3:A:466:HOH:O	1.82	0.80
1:B:107:LYS:HE3	3:B:567:HOH:O	1.81	0.79
1:A:55:LEU:HD11	1:A:187:SER:HB3	1.64	0.78
1:A:78:THR:HG22	3:A:529:HOH:O	1.84	0.77
1:A:87:ASN:O	2:A:406:IOD:I	2.73	0.76
2:A:416:IOD:I	3:A:554:HOH:O	2.74	0.74
1:A:193:LYS:O	1:A:197:VAL:HG23	1.88	0.74
1:A:108:ASN:N	1:A:151:MET:HE1	2.02	0.73
2:A:418:IOD:I	3:A:563:HOH:O	2.77	0.72
1:A:111:LYS:HG2	2:A:424:IOD:I	2.61	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:GLN:NE2	1:B:71:GLU:HG2	2.05	0.70
1:B:111:LYS:CG	2:B:419:IOD:I	3.11	0.68
2:A:433:IOD:I	3:B:476:HOH:O	2.83	0.66
1:B:107:LYS:CE	3:B:567:HOH:O	2.42	0.62
1:B:90:LYS:CE	2:B:421:IOD:I	3.17	0.61
2:B:440:IOD:I	3:B:518:HOH:O	2.87	0.57
1:A:149:ASN:C	1:A:149:ASN:HD22	2.08	0.56
1:B:86:ASP:O	1:B:87:ASN:CB	2.40	0.56
1:B:111:LYS:HE2	2:B:419:IOD:I	2.76	0.56
1:B:166:ARG:HG3	1:B:167:TRP:N	2.21	0.56
1:B:16:GLY:O	1:B:17:ILE:CB	2.56	0.53
1:B:99:LEU:HD12	1:B:103:GLU:OE2	2.07	0.53
1:A:107:LYS:C	1:A:151:MET:HE1	2.29	0.53
1:B:108:ASN:HD22	1:B:111:LYS:H	1.57	0.52
1:A:111:LYS:CG	2:A:424:IOD:I	3.27	0.52
1:B:28:LYS:HG2	1:B:57:ASP:OD2	2.09	0.52
1:B:101:GLN:HG2	3:B:513:HOH:O	2.10	0.52
1:A:49:ASN:HB2	3:A:509:HOH:O	2.09	0.51
1:A:43:VAL:HA	1:A:84:ILE:CD1	2.41	0.50
1:A:49:ASN:ND2	3:A:531:HOH:O	2.44	0.50
1:A:59:LYS:HD3	2:A:413:IOD:I	2.83	0.49
1:B:93:VAL:HG11	2:B:421:IOD:I	2.81	0.49
1:B:89:ASP:OD2	1:B:92:GLN:HG2	2.13	0.49
1:B:32:VAL:HG23	2:B:422:IOD:I	2.83	0.48
1:A:149:ASN:HD22	1:A:151:MET:H	1.62	0.46
1:A:109:PHE:HE2	2:A:402:IOD:I	2.68	0.46
1:B:72:ASP:OD2	1:B:75:ASN:HB2	2.16	0.45
1:A:107:LYS:HA	1:A:151:MET:HE2	1.98	0.45
1:B:108:ASN:HB2	2:B:414:IOD:I	2.87	0.45
1:B:27:GLY:HA3	1:B:188:LEU:HD11	1.98	0.45
1:B:148:SER:HB2	3:B:565:HOH:O	2.15	0.45
1:A:43:VAL:HA	1:A:84:ILE:HD13	1.99	0.45
1:A:109:PHE:CE2	2:A:402:IOD:I	3.41	0.44
1:B:111:LYS:CE	2:B:419:IOD:I	3.36	0.44
1:B:92:GLN:NE2	3:B:460:HOH:O	2.43	0.44
1:B:114:GLN:NE2	3:B:516:HOH:O	2.50	0.43
1:A:134:LYS:HB2	2:A:403:IOD:I	2.89	0.43
1:A:149:ASN:ND2	1:A:151:MET:H	2.16	0.43
1:B:89:ASP:O	1:B:93:VAL:HG23	2.18	0.43
1:B:107:LYS:HG3	3:B:504:HOH:O	2.17	0.43
1:B:44:ASN:HB2	2:B:426:IOD:I	2.89	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:ASN:HB3	2:A:428:IOD:I	2.89	0.43
1:B:18:LYS:HB2	1:B:18:LYS:HE2	1.39	0.43
1:A:160:VAL:HG12	1:A:162:VAL:HG13	2.01	0.42
1:A:108:ASN:HB2	2:B:415:IOD:I	2.90	0.42
1:B:19:MET:N	3:B:563:HOH:O	2.49	0.42
1:B:134:LYS:HG2	2:B:410:IOD:I	2.90	0.41
1:B:87:ASN:HA	1:B:88:GLY:O	2.20	0.41
1:B:22:THR:CG2	1:B:54:LYS:HD2	2.50	0.41
1:B:107:LYS:HD3	2:B:444:IOD:I	2.91	0.41
1:B:87:ASN:ND2	3:B:531:HOH:O	2.52	0.41
1:A:107:LYS:HA	1:A:151:MET:CE	2.51	0.40
1:B:88:GLY:O	1:B:89:ASP:C	2.58	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	178/209 (85%)	173 (97%)	5 (3%)	0	100	100
1	B	176/209 (84%)	165 (94%)	8 (4%)	3 (2%)	9	2
All	All	354/418 (85%)	338 (96%)	13 (4%)	3 (1%)	19	9

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	17	ILE
1	B	19	MET
1	B	87	ASN

### 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	148/174 (85%)	144 (97%)	4 (3%)	44	38
1	B	143/174 (82%)	134 (94%)	9 (6%)	18	8
All	All	291/348 (84%)	278 (96%)	13 (4%)	27	18

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	THR
1	A	89	ASP
1	A	149	ASN
1	A	191	LEU
1	B	18	LYS
1	B	48	PRO
1	B	49	ASN
1	B	55	LEU
1	B	72	ASP
1	B	87	ASN
1	B	90	LYS
1	B	108	ASN
1	B	166	ARG

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

Mol	Chain	Res	Type
1	A	114	GLN
1	A	149	ASN
1	A	181	HIS
1	B	92	GLN
1	B	108	ASN
1	B	114	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 monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

Of 44 ligands modelled in this entry, 44 are monoatomic - leaving 0 for Mogul analysis.

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, 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	180/209 (86%)	0.73	12 (6%) 17 20	24, 30, 38, 42	0
1	B	178/209 (85%)	0.94	24 (13%) 3 3	24, 30, 46, 62	0
All	All	358/418 (85%)	0.84	36 (10%) 7 8	24, 30, 40, 62	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	189	LEU	6.4
1	B	188	LEU	5.6
1	B	16	GLY	5.6
1	B	59	LYS	5.5
1	B	19	MET	5.2
1	B	17	ILE	5.0
1	A	198	GLY	4.9
1	A	197	VAL	4.7
1	B	191	LEU	4.5
1	A	19	MET	4.3
1	B	20	PRO	4.1
1	B	185	SER	4.1
1	B	125	LEU	3.8
1	B	173	MET	3.7
1	A	133	LEU	3.6
1	B	87	ASN	3.3
1	B	190	GLU	3.2
1	B	88	GLY	3.0
1	B	57	ASP	2.8
1	B	186	LYS	2.8
1	A	104	LEU	2.7
1	A	88	GLY	2.7
1	B	187	SER	2.7
1	A	89	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	61	ALA	2.5
1	B	133	LEU	2.5
1	B	18	LYS	2.5
1	A	196	ASN	2.4
1	A	82	ALA	2.3
1	A	68	ASP	2.2
1	B	28	LYS	2.2
1	A	99	LEU	2.2
1	B	192	LYS	2.2
1	B	105	VAL	2.2
1	A	79	SER	2.1
1	B	193	LYS	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(Å <sup>2</sup> )	Q<0.9
2	IOD	A	434	1/1	0.77	0.16	78,78,78,78	1
2	IOD	B	441	1/1	0.79	0.11	48,48,48,48	1
2	IOD	A	402	1/1	0.84	0.17	43,43,43,43	1
2	IOD	B	405	1/1	0.87	0.11	60,60,60,60	1
2	IOD	B	435	1/1	0.90	0.11	69,69,69,69	1
2	IOD	A	445	1/1	0.90	0.08	63,63,63,63	1
2	IOD	A	431	1/1	0.91	0.11	60,60,60,60	1
2	IOD	A	442	1/1	0.92	0.13	64,64,64,64	1
2	IOD	B	420	1/1	0.92	0.09	55,55,55,55	1
2	IOD	B	438	1/1	0.93	0.06	52,52,52,52	1
2	IOD	B	421	1/1	0.93	0.12	53,53,53,53	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	IOD	B	444	1/1	0.93	0.08	65,65,65,65	1
2	IOD	A	433	1/1	0.94	0.14	45,45,45,45	1
2	IOD	B	440	1/1	0.94	0.10	48,48,48,48	1
2	IOD	B	429	1/1	0.94	0.12	62,62,62,62	1
2	IOD	A	428	1/1	0.94	0.17	43,43,43,43	1
2	IOD	B	426	1/1	0.95	0.09	36,36,36,36	1
2	IOD	A	436	1/1	0.95	0.09	53,53,53,53	1
2	IOD	A	408	1/1	0.95	0.17	36,36,36,36	1
2	IOD	B	437	1/1	0.95	0.09	58,58,58,58	1
2	IOD	A	427	1/1	0.96	0.12	52,52,52,52	1
2	IOD	B	443	1/1	0.96	0.10	44,44,44,44	1
2	IOD	B	432	1/1	0.96	0.07	59,59,59,59	1
2	IOD	B	415	1/1	0.97	0.08	43,43,43,43	1
2	IOD	B	419	1/1	0.97	0.07	61,61,61,61	1
2	IOD	A	411	1/1	0.97	0.07	47,47,47,47	1
2	IOD	A	416	1/1	0.97	0.08	49,49,49,49	1
2	IOD	B	439	1/1	0.97	0.05	60,60,60,60	1
2	IOD	B	422	1/1	0.97	0.07	42,42,42,42	1
2	IOD	B	423	1/1	0.97	0.13	41,41,41,41	1
2	IOD	A	424	1/1	0.97	0.14	43,43,43,43	1
2	IOD	B	409	1/1	0.97	0.10	48,48,48,48	1
2	IOD	A	401	1/1	0.98	0.09	42,42,42,42	1
2	IOD	A	406	1/1	0.98	0.06	36,36,36,36	1
2	IOD	B	410	1/1	0.98	0.08	49,49,49,49	1
2	IOD	A	430	1/1	0.98	0.08	74,74,74,74	1
2	IOD	A	418	1/1	0.99	0.11	34,34,34,34	1
2	IOD	A	412	1/1	0.99	0.04	41,41,41,41	1
2	IOD	B	414	1/1	0.99	0.05	44,44,44,44	1
2	IOD	A	413	1/1	0.99	0.13	36,36,36,36	1
2	IOD	B	417	1/1	0.99	0.15	35,35,35,35	1
2	IOD	A	403	1/1	0.99	0.10	39,39,39,39	1
2	IOD	B	407	1/1	0.99	0.07	32,32,32,32	1
2	IOD	B	404	1/1	1.00	0.15	28,28,28,28	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.