



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2018 – 12:46 PM EST

PDB ID : 5Y7Y
Title : Crystal structure of AhRR/ARNT complex
Authors : Sakurai, S.; Shimizu, T.; Ohto, U.
Deposited on : 2017-08-18
Resolution : 2.40 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030736
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030736

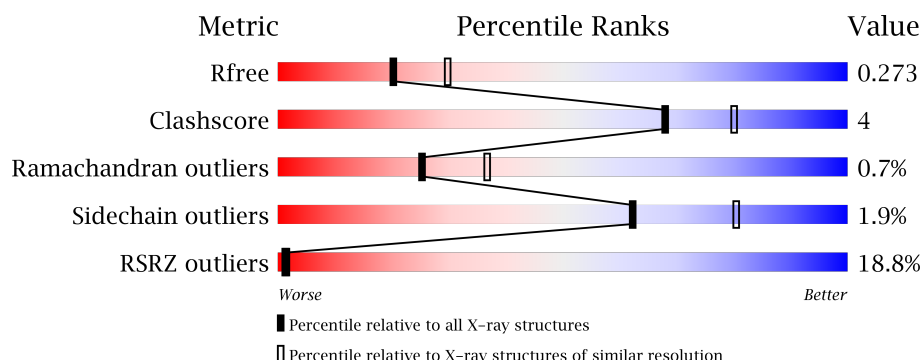
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	100719	3166 (2.40-2.40)
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)
RSRZ outliers	101464	3195 (2.40-2.40)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	258	<div> <div>2%</div> <div>73%</div> <div>5%</div> <div>22%</div> </div>
2	B	311	<div> <div>26%</div> <div>75%</div> <div>9%</div> <div>15%</div> </div>

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
3	GOL	B	501	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3782 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 Aryl hydrocarbon receptor repressor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	201	Total	C	N	O	S	0	0	0
			1615	1040	280	286	9			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	GLY	-	expression tag	UNP A9YTQ3
A	24	PRO	-	expression tag	UNP A9YTQ3
A	25	GLU	-	expression tag	UNP A9YTQ3
A	26	PHE	-	expression tag	UNP A9YTQ3

- Molecule 2 is a protein called Aryl hydrocarbon receptor nuclear translocator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	263	Total	C	N	O	S	0	0	0
			2127	1340	379	393	15			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	78	GLY	-	expression tag	UNP Q9BE97
B	79	PRO	-	expression tag	UNP Q9BE97
B	80	GLU	-	expression tag	UNP Q9BE97
B	81	PHE	-	expression tag	UNP Q9BE97
B	?	-	THR	deletion	UNP Q9BE97
B	?	-	SER	deletion	UNP Q9BE97
B	?	-	THR	deletion	UNP Q9BE97
B	?	-	ASP	deletion	UNP Q9BE97
B	?	-	GLY	deletion	UNP Q9BE97
B	?	-	THR	deletion	UNP Q9BE97
B	?	-	TYR	deletion	UNP Q9BE97
B	?	-	ALA	deletion	UNP Q9BE97

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	LEU	deletion	UNP Q9BE97
B	?	-	THR	deletion	UNP Q9BE97
B	?	-	GLY	deletion	UNP Q9BE97
B	?	-	ARG	deletion	UNP Q9BE97
B	?	-	ILE	deletion	UNP Q9BE97
B	?	-	LEU	deletion	UNP Q9BE97
B	?	-	ASP	deletion	UNP Q9BE97
B	?	-	LEU	deletion	UNP Q9BE97
B	?	-	LYS	deletion	UNP Q9BE97
B	?	-	THR	deletion	UNP Q9BE97
B	?	-	GLY	deletion	UNP Q9BE97
B	?	-	THR	deletion	UNP Q9BE97
B	?	-	VAL	deletion	UNP Q9BE97
B	?	-	LYS	deletion	UNP Q9BE97
B	?	-	LYS	deletion	UNP Q9BE97
B	?	-	GLU	deletion	UNP Q9BE97
B	?	-	GLY	deletion	UNP Q9BE97
B	?	-	GLN	deletion	UNP Q9BE97
B	?	-	GLN	deletion	UNP Q9BE97
B	?	-	SER	deletion	UNP Q9BE97
B	?	-	SER	deletion	UNP Q9BE97
B	?	-	MET	deletion	UNP Q9BE97
B	?	-	ARG	deletion	UNP Q9BE97
B	?	-	MET	deletion	UNP Q9BE97
B	?	-	CYS	deletion	UNP Q9BE97
B	?	-	SER	deletion	UNP Q9BE97
B	?	-	SER	deletion	UNP Q9BE97
B	?	-	VAL	deletion	UNP Q9BE97
B	?	-	ASP	deletion	UNP Q9BE97
B	?	-	SER	deletion	UNP Q9BE97
B	?	-	VAL	deletion	UNP Q9BE97
B	?	-	SER	deletion	UNP Q9BE97
B	?	-	MET	deletion	UNP Q9BE97
B	?	-	ASN	deletion	UNP Q9BE97
B	?	-	ARG	deletion	UNP Q9BE97
B	?	-	LEU	deletion	UNP Q9BE97
B	?	-	SER	deletion	UNP Q9BE97
B	?	-	PHE	deletion	UNP Q9BE97
B	?	-	VAL	deletion	UNP Q9BE97
B	?	-	ARG	deletion	UNP Q9BE97
B	?	-	ASN	deletion	UNP Q9BE97
B	?	-	ARG	deletion	UNP Q9BE97

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	CYS	deletion	UNP Q9BE97
B	?	-	ARG	deletion	UNP Q9BE97
B	?	-	ASN	deletion	UNP Q9BE97
B	?	-	GLY	deletion	UNP Q9BE97
B	?	-	LEU	deletion	UNP Q9BE97
B	?	-	GLY	deletion	UNP Q9BE97
B	?	-	SER	deletion	UNP Q9BE97
B	?	-	ALA	deletion	UNP Q9BE97
B	?	-	LYS	deletion	UNP Q9BE97
B	?	-	ASP	deletion	UNP Q9BE97
B	?	-	GLY	deletion	UNP Q9BE97
B	?	-	GLU	deletion	UNP Q9BE97
B	?	-	ALA	deletion	UNP Q9BE97
B	?	-	GLY	deletion	UNP Q9BE97
B	?	-	VAL	deletion	UNP Q9BE97
B	?	-	SER	deletion	UNP Q9BE97
B	?	-	LEU	deletion	UNP Q9BE97
B	?	-	PRO	deletion	UNP Q9BE97
B	?	-	ASP	deletion	UNP Q9BE97
B	?	-	ASP	deletion	UNP Q9BE97
B	?	-	ASP	deletion	UNP Q9BE97
B	?	-	PRO	deletion	UNP Q9BE97
B	?	-	GLU	deletion	UNP Q9BE97
B	?	-	ALA	deletion	UNP Q9BE97
B	?	-	GLY	deletion	UNP Q9BE97
B	?	-	GLN	deletion	UNP Q9BE97

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

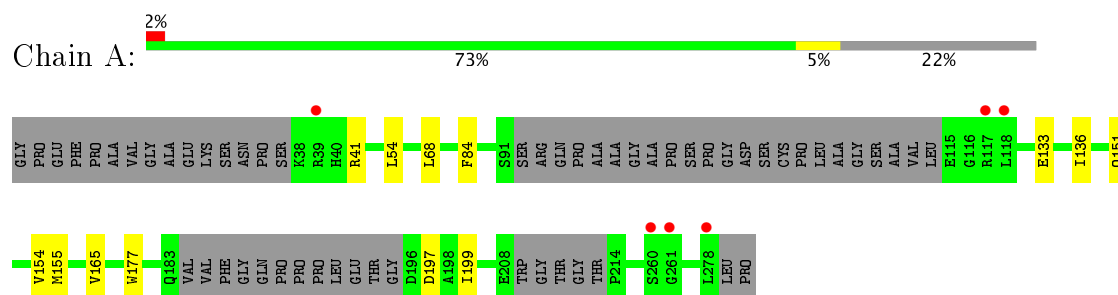
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	14	Total	O	0	0
			14	14		
4	B	8	Total	O	0	0
			8	8		

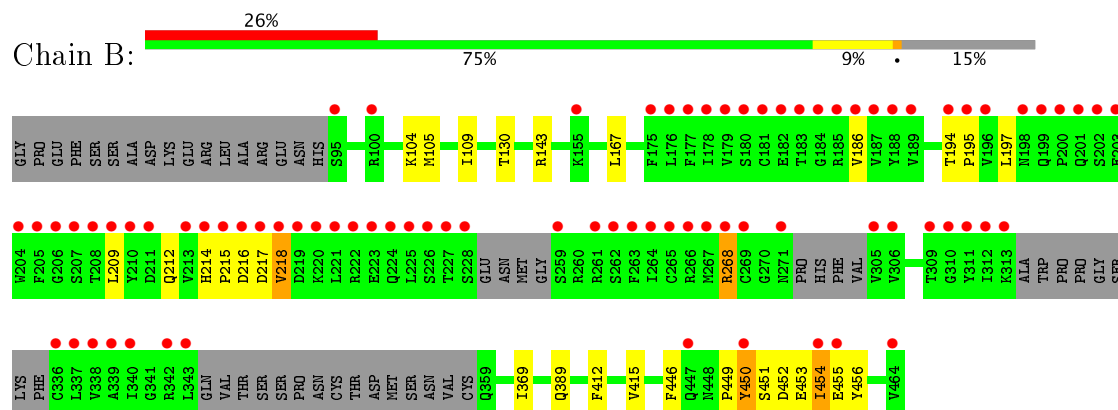
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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: Aryl hydrocarbon receptor repressor



- Molecule 2: Aryl hydrocarbon receptor nuclear translocator



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	78.36 Å 78.36 Å 129.78 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 49.98 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.00-2.40) 99.9 (49.98-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.95 (at 2.39 Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.234 , 0.268 0.237 , 0.273	Depositor DCC
R_{free} test set	1542 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	54.9	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 46.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.047 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3782	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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

5.1 Standard geometry

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	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.30	0/1649	0.49	0/2224
2	B	0.32	0/2166	0.54	0/2920
All	All	0.31	0/3815	0.52	0/5144

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

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	1615	0	1629	8	3
2	B	2127	0	2118	28	3
3	A	12	0	16	0	0
3	B	6	0	8	0	0
4	A	14	0	0	1	0
4	B	8	0	0	0	0
All	All	3782	0	3771	33	3

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 (33) 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:369:ILE:HD11	2:B:454:ILE:HG12	1.44	0.98
2:B:454:ILE:HD13	2:B:454:ILE:H	1.46	0.81
1:A:136:ILE:HD13	1:A:154:VAL:HG13	1.75	0.69
2:B:217:ASP:HA	2:B:218:VAL:HB	1.74	0.69
2:B:369:ILE:CD1	2:B:454:ILE:HG12	2.20	0.69
2:B:369:ILE:HD12	2:B:369:ILE:H	1.64	0.61
1:A:151:GLN:O	1:A:155:MET:HG2	2.02	0.58
2:B:369:ILE:HD12	2:B:369:ILE:N	2.23	0.54
2:B:451:SER:HB2	2:B:453:GLU:HB2	1.90	0.54
2:B:369:ILE:CD1	2:B:454:ILE:CG1	2.85	0.53
1:A:177:TRP:CZ2	1:A:199:ILE:HG21	2.44	0.51
2:B:446:PHE:CZ	2:B:456:TYR:HD2	2.28	0.51
2:B:451:SER:C	2:B:453:GLU:N	2.65	0.50
2:B:455:GLU:OE1	2:B:455:GLU:N	2.44	0.50
1:A:68:LEU:HD11	2:B:104:LYS:HB3	1.95	0.49
2:B:451:SER:CB	2:B:453:GLU:HB2	2.42	0.49
2:B:449:PRO:O	2:B:450:TYR:CB	2.60	0.49
2:B:453:GLU:OE2	2:B:453:GLU:HA	2.14	0.48
1:A:165:VAL:HG23	4:A:413:HOH:O	2.14	0.47
2:B:212:GLN:HE22	2:B:268:ARG:HE	1.64	0.46
2:B:369:ILE:HD11	2:B:454:ILE:CG1	2.28	0.46
1:A:136:ILE:HG21	1:A:154:VAL:CG1	2.46	0.45
2:B:455:GLU:CD	2:B:455:GLU:H	2.19	0.45
2:B:214:HIS:HB2	2:B:215:PRO:HA	2.00	0.44
2:B:454:ILE:HB	2:B:455:GLU:OE1	2.18	0.44
2:B:451:SER:OG	2:B:453:GLU:HB2	2.18	0.44
2:B:105:MET:O	2:B:109:ILE:HG12	2.19	0.43
1:A:54:LEU:O	2:B:143:ARG:HD3	2.18	0.43
2:B:412:PHE:HA	2:B:415:VAL:HG22	2.01	0.42
2:B:194:THR:N	2:B:195:PRO:HD2	2.35	0.42
1:A:84:PHE:HZ	2:B:167:LEU:HD13	1.85	0.41
2:B:446:PHE:CZ	2:B:456:TYR:CD2	3.09	0.41
2:B:215:PRO:HA	2:B:216:ASP:HA	1.84	0.41

All (3) 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:41:ARG:CZ	2:B:452:ASP:OD1[3_555]	1.56	0.64
1:A:41:ARG:NE	2:B:452:ASP:OD1[3_555]	1.84	0.36

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:ARG:NH2	2:B:452:ASP:OD1[3_555]	2.03	0.17

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	193/258 (75%)	191 (99%)	2 (1%)	0	100	100
2	B	253/311 (81%)	240 (95%)	10 (4%)	3 (1%)	15	21
All	All	446/569 (78%)	431 (97%)	12 (3%)	3 (1%)	25	37

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	218	VAL
2	B	450	TYR
2	B	186	VAL

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	178/219 (81%)	176 (99%)	2 (1%)	78	90
2	B	243/285 (85%)	237 (98%)	6 (2%)	53	73
All	All	421/504 (84%)	413 (98%)	8 (2%)	62	80

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	133	GLU
1	A	197	ASP
2	B	130	THR
2	B	197	LEU
2	B	209	LEU
2	B	268	ARG
2	B	389	GLN
2	B	454	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	176	HIS
1	A	243	GLN
2	B	103	ASN
2	B	212	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

3 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	$\# Z > 2$
3	GOL	A	301	-	5,5,5	0.20	0	5,5,5	0.27	0
3	GOL	A	302	-	5,5,5	0.19	0	5,5,5	0.29	0
3	GOL	B	501	-	5,5,5	0.22	0	5,5,5	0.30	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	GOL	A	301	-	-	0/4/4/4	0/0/0/0
3	GOL	A	302	-	-	0/4/4/4	0/0/0/0
3	GOL	B	501	-	-	0/4/4/4	0/0/0/0

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.

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	201/258 (77%)	0.27	6 (2%)	51	49	37, 51, 99, 125	0
2	B	263/311 (84%)	1.86	81 (30%)	0	0	40, 76, 172, 191	0
All	All	464/569 (81%)	1.17	87 (18%)	1	1	37, 59, 167, 191	0

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	210	TYR	12.1
2	B	179	VAL	10.5
2	B	336	CYS	9.4
2	B	311	TYR	9.2
2	B	208	THR	8.5
2	B	204	TRP	8.1
2	B	188	TYR	8.0
2	B	337	LEU	7.9
2	B	189	VAL	7.9
2	B	177	PHE	7.7
2	B	216	ASP	7.6
2	B	183	THR	7.5
2	B	213	VAL	7.5
2	B	215	PRO	7.4
2	B	313	LYS	7.4
2	B	225	LEU	7.3
2	B	181	CYS	7.3
2	B	221	LEU	7.2
2	B	185	ARG	6.9
2	B	178	ILE	6.6
2	B	450	TYR	6.5
2	B	219	ASP	6.3
2	B	267	MET	5.9
2	B	214	HIS	5.7

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Mol	Chain	Res	Type	RSRZ
2	B	211	ASP	5.7
2	B	187	VAL	5.6
2	B	184	GLY	5.6
2	B	182	GLU	5.5
2	B	217	ASP	5.5
2	B	202	SER	5.5
2	B	222	ARG	5.4
2	B	209	LEU	5.4
2	B	186	VAL	5.3
2	B	310	GLY	5.2
2	B	224	GLN	5.1
2	B	338	VAL	5.1
2	B	223	GLU	5.1
2	B	306	VAL	5.0
2	B	227	THR	4.9
2	B	226	SER	4.9
2	B	207	SER	4.8
2	B	95	SER	4.8
2	B	269	CYS	4.6
2	B	454	ILE	4.6
2	B	305	VAL	4.6
2	B	205	PHE	4.5
2	B	218	VAL	4.5
2	B	203	GLU	4.4
2	B	312	ILE	4.3
2	B	200	PRO	4.3
2	B	175	PHE	4.2
2	B	263	PHE	4.1
2	B	343	LEU	4.1
2	B	199	GLN	4.1
2	B	265	CYS	4.1
2	B	194	THR	4.1
2	B	196	VAL	3.9
2	B	261	ARG	3.8
2	B	455	GLU	3.8
2	B	264	ILE	3.7
2	B	266	ARG	3.7
2	B	268	ARG	3.7
2	B	180	SER	3.6
2	B	228	SER	3.5
2	B	206	GLY	3.4
2	B	220	LYS	3.4

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Mol	Chain	Res	Type	RSRZ
2	B	198	ASN	3.2
2	B	259	SER	3.2
2	B	176	LEU	3.1
1	A	260	SER	3.1
1	A	117	ARG	3.0
2	B	262	SER	2.9
2	B	201	GLN	2.8
2	B	342	ARG	2.8
2	B	100	ARG	2.7
2	B	464	VAL	2.7
2	B	195	PRO	2.6
1	A	39	ARG	2.5
1	A	278	LEU	2.5
1	A	118	LEU	2.4
2	B	447	GLN	2.3
2	B	340	ILE	2.3
2	B	339	ALA	2.3
2	B	309	THR	2.3
2	B	271	ASN	2.3
1	A	261	GLY	2.1
2	B	155	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	GOL	B	501	6/6	0.82	0.29	5.00	72,74,76,77	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	GOL	A	301	6/6	0.94	0.23	1.90	60,62,62,65	0
3	GOL	A	302	6/6	0.93	0.19	0.80	73,73,74,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.