



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:53 PM GMT

PDB ID : 4GJQ
Title : Crystal structure of human GLTP bound with 12:0 monosulfatide (orthorhombic form; two subunits in asymmetric unit)
Authors : Cabo-Bilbao, A.; Goni-de-Cerio, F.; Samygina, V.R.; Ochoa-Lizarralde, B.; Popov, A.N.; Malinina, L.
Deposited on : 2012-08-10
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
<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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

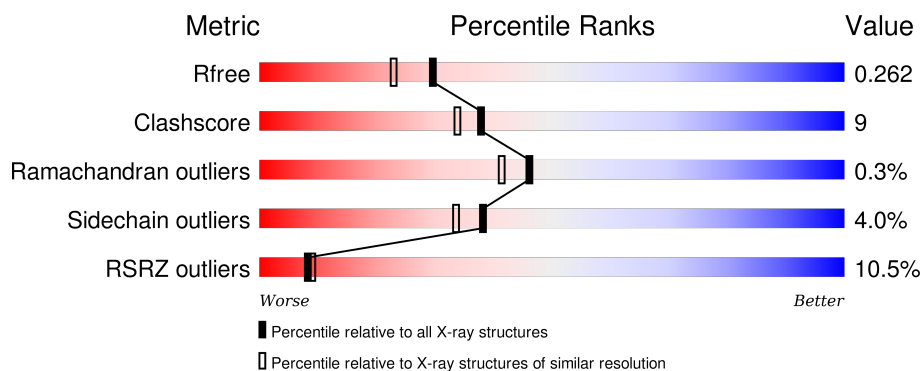
1 Overall quality at a glance

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}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (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 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	209	
1	B	209	

2 Entry composition [i](#)

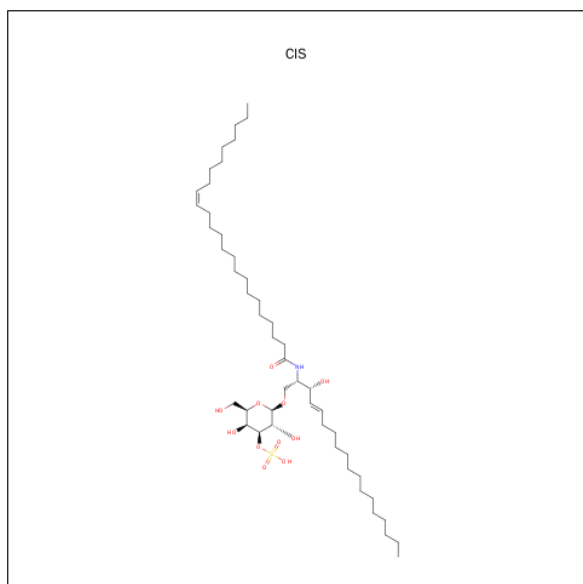
There are 4 unique types of molecules in this entry. The entry contains 3599 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glycolipid transfer protein.

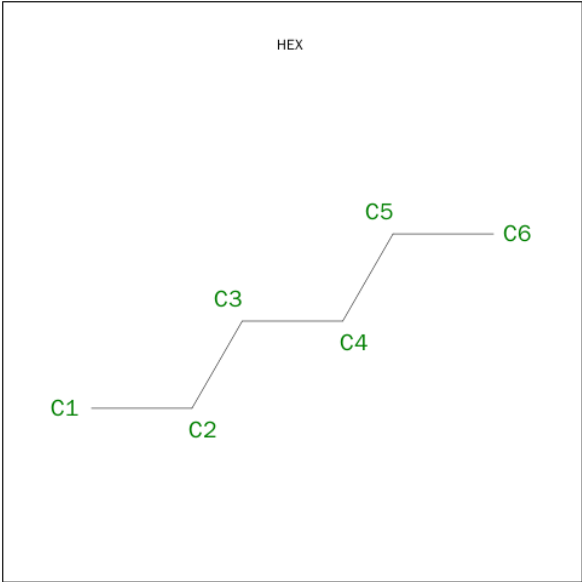
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	198	Total	C	N	O	S	0	1	0
			1598	1044	258	288	8			
1	B	202	Total	C	N	O	S	0	3	0
			1630	1061	265	296	8			

- Molecule 2 is (15Z)-N-((1S,2R,3E)-2-HYDROXY-1-[(3-O-SULFO-BETA-D-GALACTOPYRANOSYL)OXY]METHYL)HEPTADEC-3-ENYL)TETRACOS-15-ENAMIDE (three-letter code: CIS) (formula: C₄₈H₉₁NO₁₁S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			49	36	1	11	1		
2	B	1	Total	C	N	O	S	0	0
			49	36	1	11	1		

- Molecule 3 is HEXANE (three-letter code: HEX) (formula: C₆H₁₄).



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

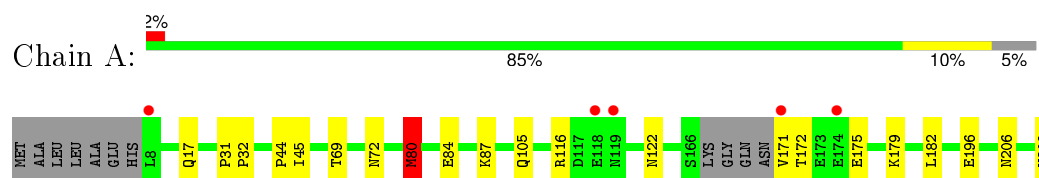
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	135	Total O 135 135	0	0
4	B	132	Total O 132 132	0	0

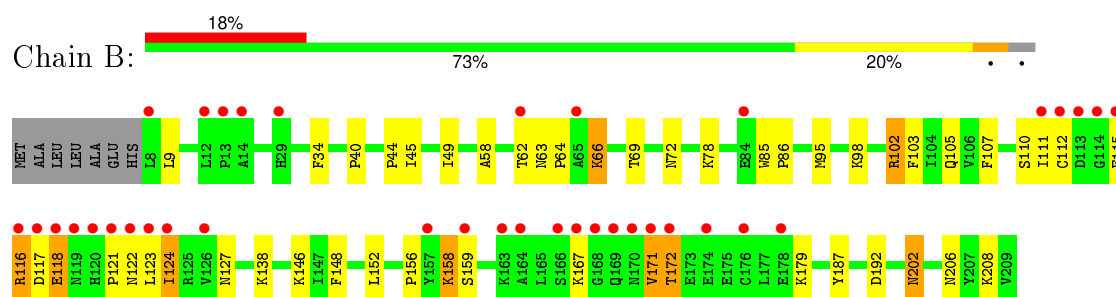
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Glycolipid transfer protein



• Molecule 1: Glycolipid transfer protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	50.96 Å 95.47 Å 132.39 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.00 15.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.3 (15.00-2.00) 98.3 (15.00-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 2.00 Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.217 , 0.260 0.220 , 0.262	Depositor DCC
R_{free} test set	2250 reflections (5.43%)	DCC
Wilson B-factor (Å ²)	29.6	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 55.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 43616 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3599	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.08% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: CIS, HEX

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	0.88	0/1636	0.81	2/2216 (0.1%)
1	B	0.89	0/1679	0.84	3/2274 (0.1%)
All	All	0.89	0/3315	0.82	5/4490 (0.1%)

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	1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	124	ILE	CB-CA-C	-8.09	95.43	111.60
1	A	80	MET	CA-CB-CG	6.23	123.89	113.30
1	B	102	ARG	NE-CZ-NH2	6.21	123.41	120.30
1	B	102	ARG	NE-CZ-NH1	-5.59	117.50	120.30
1	A	80	MET	CB-CA-C	5.33	121.06	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	118	GLU	Peptide

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	1598	0	1614	17	0
1	B	1630	0	1653	47	0
2	A	49	0	65	4	0
2	B	49	0	65	2	0
3	A	6	0	14	2	0
4	A	135	0	0	8	0
4	B	132	0	0	10	0
All	All	3599	0	3411	62	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 9.

All (62) 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:B:124:ILE:O	1:B:124:ILE:HG22	1.57	1.00
1:B:115:GLU:O	1:B:116:ARG:HB3	1.73	0.89
1:B:102:ARG:HH11	1:B:105:GLN:HE22	1.20	0.88
1:A:172:THR:HG23	4:A:456:HOH:O	1.76	0.86
1:B:158:LYS:HE2	1:B:159:SER:H	1.40	0.85
2:A:301:CIS:H142	1:B:45:ILE:HD11	1.60	0.81
1:A:87:LYS:NZ	4:A:530:HOH:O	1.82	0.81
1:B:156:PRO:O	4:B:916:HOH:O	2.05	0.75
1:B:124:ILE:CG2	1:B:124:ILE:O	2.29	0.72
1:B:118:GLU:HG2	1:B:121:PRO:HB3	1.74	0.70
1:A:80:MET:HE1	4:A:505:HOH:O	1.91	0.69
1:A:206:ASN:H	3:A:302:HEX:H32	1.58	0.67
1:B:124:ILE:HG23	1:B:127:ASN:HD22	1.62	0.65
1:B:102:ARG:HH11	1:B:105:GLN:NE2	1.94	0.64
1:B:158:LYS:HE2	1:B:159:SER:N	2.14	0.61
1:B:102:ARG:NH1	1:B:192:ASP:OD1	2.34	0.61
1:B:124:ILE:CG2	1:B:127:ASN:HB2	2.32	0.60
1:B:116:ARG:HG2	1:B:117:ASP:N	2.17	0.59
1:A:44:PRO:HG2	2:B:700:CIS:H91	1.84	0.58
1:B:107:PHE:CE1	1:B:124:ILE:HG21	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:208:LYS:NZ	4:B:914:HOH:O	2.36	0.56
1:B:34:PHE:CE2	1:B:49:ILE:HD12	2.40	0.56
1:B:118:GLU:HA	1:B:121:PRO:HA	1.88	0.55
1:A:206:ASN:N	3:A:302:HEX:H32	2.23	0.54
1:B:78:LYS:HG3	4:B:932:HOH:O	2.09	0.53
1:B:202:ASN:HB2	4:B:870:HOH:O	2.08	0.52
1:B:116:ARG:HD2	1:B:118:GLU:OE2	2.09	0.52
2:A:301:CIS:H151	1:B:148:PHE:HA	1.91	0.52
1:B:146:LYS:HG3	4:B:926:HOH:O	2.09	0.52
1:B:58:ALA:O	1:B:62:THR:HG23	2.10	0.52
1:B:9:LEU:HD22	1:B:187:TYR:HA	1.92	0.51
1:A:45:ILE:HD13	2:A:301:CIS:H251	1.92	0.51
1:A:209:VAL:HG23	1:B:40:PRO:HG2	1.92	0.50
1:B:85:TRP:CG	1:B:86:PRO:HA	2.48	0.49
1:B:66:LYS:NZ	4:B:838:HOH:O	2.13	0.48
1:B:102:ARG:NH1	1:B:105:GLN:HE22	1.99	0.48
1:B:171:VAL:HG23	1:B:172:THR:O	2.14	0.47
1:A:80:MET:CE	4:A:505:HOH:O	2.58	0.46
1:A:122:ASN:HB3	4:A:495:HOH:O	2.14	0.46
1:A:172:THR:O	1:A:175:GLU:HG3	2.16	0.46
2:A:301:CIS:H91	1:B:44:PRO:HG2	1.98	0.46
1:B:123:LEU:HD12	4:B:890:HOH:O	2.16	0.46
1:B:111:ILE:HG22	1:B:112:CYS:N	2.31	0.45
1:A:17:GLN:NE2	4:A:431:HOH:O	2.40	0.45
1:A:69:THR:H	1:A:72:ASN:HD22	1.65	0.44
1:B:206:ASN:ND2	4:B:815:HOH:O	2.43	0.44
1:B:103:PHE:CE1	2:B:700:CIS:H252	2.53	0.43
1:B:110:SER:HA	4:B:835:HOH:O	2.18	0.43
1:B:122:ASN:HB2	1:B:158:LYS:CG	2.49	0.43
1:B:138:LYS:HE2	1:B:206:ASN:ND2	2.33	0.43
1:A:31:PRO:HB2	1:A:32:PRO:HD3	2.01	0.43
1:B:63:ASN:O	1:B:64:PRO:C	2.57	0.42
1:A:209:VAL:HG23	1:B:40:PRO:CG	2.49	0.42
1:B:124:ILE:HG22	1:B:127:ASN:HB2	2.02	0.42
1:A:84:GLU:HG2	4:A:513:HOH:O	2.19	0.41
1:B:118:GLU:HA	1:B:121:PRO:CA	2.51	0.41
1:B:34:PHE:HE2	1:B:49:ILE:HD12	1.82	0.41
1:B:124:ILE:HG23	1:B:127:ASN:ND2	2.29	0.41
1:B:123:LEU:O	4:B:849:HOH:O	2.22	0.41
1:B:69:THR:HG23	1:B:72:ASN:ND2	2.36	0.41
1:A:105:GLN:NE2	4:A:442:HOH:O	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:MET:O	1:B:98:LYS:HG2	2.22	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	194/209 (93%)	190 (98%)	4 (2%)	0	100	100
1	B	202/209 (97%)	189 (94%)	12 (6%)	1 (0%)	34	26
All	All	396/418 (95%)	379 (96%)	16 (4%)	1 (0%)	46	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	167	LYS

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	172/180 (96%)	166 (96%)	6 (4%)	43	40
1	B	177/180 (98%)	169 (96%)	8 (4%)	34	29
All	All	349/360 (97%)	335 (96%)	14 (4%)	38	33

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	MET
1	A	116	ARG
1	A	171	VAL
1	A	179	LYS
1	A	182	LEU
1	A	196	GLU
1	B	66	LYS
1	B	116	ARG
1	B	152	LEU
1	B	158	LYS
1	B	171	VAL
1	B	172	THR
1	B	179	LYS
1	B	202	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	72	ASN
1	A	120	HIS
1	A	127	ASN
1	B	105	GLN
1	B	120	HIS
1	B	127	ASN
1	B	186	ASN
1	B	202	ASN
1	B	206	ASN

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 [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$
2	CIS	A	301	-	48,49,61	0.71	1 (2%)	54,60,72	1.00	2 (3%)
3	HEX	A	302	-	5,5,5	0.39	0	4,4,4	0.05	0
2	CIS	B	700	-	48,49,61	0.69	1 (2%)	54,60,72	1.11	5 (9%)

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
2	CIS	A	301	-	-	0/46/66/78	0/1/1/1
3	HEX	A	302	-	-	0/3/3/3	0/0/0/0
2	CIS	B	700	-	-	0/46/66/78	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	CIS	O-C43	-2.31	1.36	1.40
2	B	700	CIS	O5-C46	2.16	1.48	1.43

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	700	CIS	O2-C18-C19	-3.49	115.95	121.98
2	B	700	CIS	C1-N-C18	-3.01	118.04	123.61
2	B	700	CIS	C-C1-N	-2.30	106.55	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	CIS	O2-C18-C19	-2.02	118.51	121.98
2	B	700	CIS	C19-C18-N	2.14	119.32	115.83
2	A	301	CIS	C45-O4-S	2.28	123.11	118.77
2	B	700	CIS	O4-C45-C46	2.35	113.64	108.43

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	CIS	4	0
3	A	302	HEX	2	0
2	B	700	CIS	2	0

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	198/209 (94%)	-0.14	5 (2%) 61 61	19, 36, 67, 84	1 (0%)
1	B	202/209 (96%)	0.68	37 (18%) 2 2	20, 47, 103, 126	0
All	All	400/418 (95%)	0.27	42 (10%) 8 9	19, 39, 91, 126	1 (0%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	168	GLY	6.8
1	B	170	ASN	5.1
1	B	167	LYS	5.0
1	B	14	ALA	4.5
1	A	174	GLU	4.5
1	B	65	ALA	4.3
1	B	122	ASN	4.2
1	B	157	TYR	4.2
1	B	169	GLN	4.1
1	B	178	GLU	3.7
1	B	163	LYS	3.6
1	B	113	ASP	3.3
1	B	121	PRO	3.3
1	B	8	LEU	3.3
1	B	62	THR	3.3
1	B	111	ILE	3.1
1	B	13	PRO	3.0
1	B	112	CYS	2.9
1	B	84	GLU	2.9
1	B	118	GLU	2.8
1	B	114	GLY	2.8
1	A	8	LEU	2.7
1	B	12	LEU	2.7
1	B	123	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	119	ASN	2.6
1	B	159	SER	2.6
1	B	166	SER	2.6
1	B	172	THR	2.5
1	B	164	ALA	2.4
1	B	115	GLU	2.4
1	B	29	HIS	2.3
1	A	171	VAL	2.3
1	B	117	ASP	2.3
1	B	174	GLU	2.3
1	B	176	CYS	2.3
1	B	124	ILE	2.3
1	A	118	GLU	2.3
1	B	116	ARG	2.2
1	A	119	ASN	2.2
1	B	171	VAL	2.1
1	B	120	HIS	2.0
1	B	126	VAL	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
2	CIS	A	301	49/61	0.96	0.10	0.11	18,24,43,46	0
2	CIS	B	700	49/61	0.97	0.09	-0.56	18,22,34,38	0
3	HEX	A	302	6/6	0.77	0.24	-	39,44,47,51	0

6.5 Other polymers [i](#)

There are no such residues in this entry.