



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:45 PM UTC

PDB ID : 4WWK / pdb_00004wwk
Title : Crystal structure of human TCR Alpha Chain-TRAV12-3, Beta Chain-TRBV6-5, Antigen-presenting molecule CD1d, and Beta-2-microglobulin
Authors : Le Nours, J.; Praveena, T.; Pellicci, D.G.; Gherardin, N.A.; Lim, R.T.; Besra, G.; Keshipeddy, A.; Richardson, S.K.; Howell, A.R.; Gras, S.; Godfrey, D.I.; Rossjohn, J.; Uldrich, A.P.
Deposited on : 2014-11-11
Resolution : 3.10 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 ⓘ 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:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

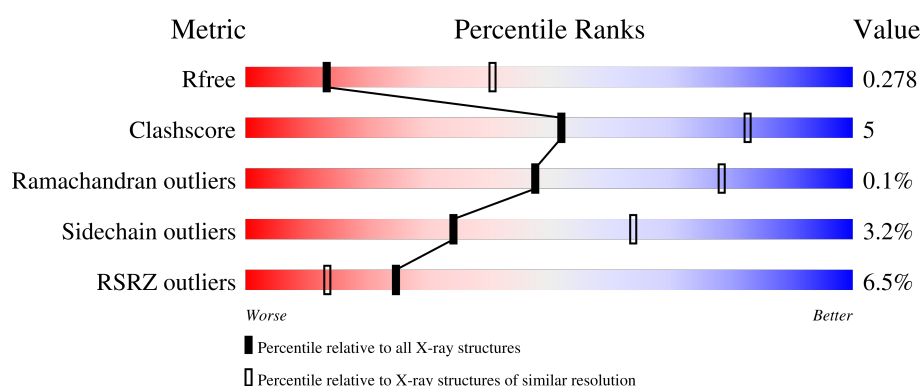
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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 $\leq 5\%$. 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	12% 81% 8% 10%
2	B	243	8% 89% 9%
3	C	278	% 81% 16%
4	D	119	3% 71% 12% 17%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 5883 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 TCR Alpha Chain-TRAV12-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	188	Total	C	N	O	S	0	0	0
			1265	780	214	263	8			

- Molecule 2 is a protein called TCR Beta Chain-TRBV6-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	240	Total	C	N	O	S	0	0	0
			1658	1042	288	320	8			

- Molecule 3 is a protein called Antigen-presenting glycoprotein CD1d.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	270	Total	C	N	O	S	0	0	0
			2088	1339	359	383	7			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	SER	-	expression tag	UNP P15813
C	1	PRO	-	expression tag	UNP P15813
C	2	GLY	-	expression tag	UNP P15813

- Molecule 4 is a protein called Beta-2-microglobulin.

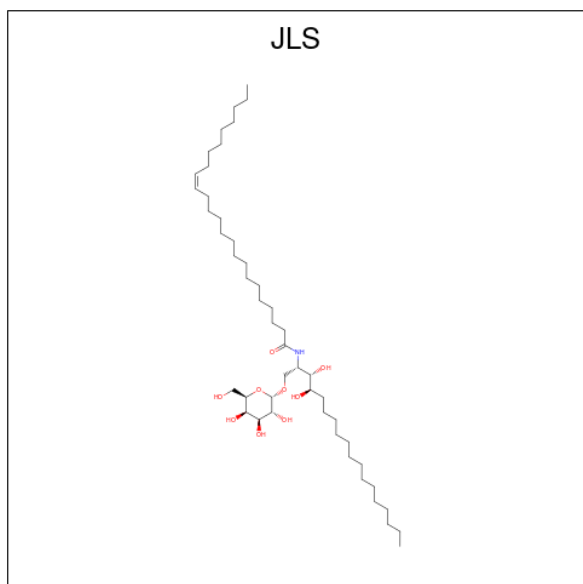
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	99	Total	C	N	O	S	0	0	0
			783	501	132	147	3			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	C	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		

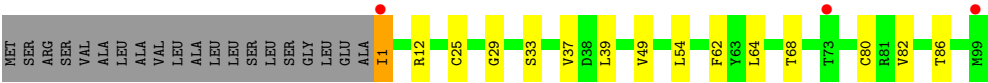
- Molecule 6 is (15Z)-N-[(2S,3S,4R)-1-(alpha-D-galactopyranosyloxy)-3,4-dihydroxyoctadecan-2-yl]tetracos-15-enamide (CCD ID: JLS) (formula: C₄₈H₉₃NO₉).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	C	1	Total	C	N	O	0	0
			58	48	1	9		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	2	Total 2	O 2	0	0
7	D	1	Total 1	O 1	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	56.38Å 76.07Å 255.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.61 – 3.10 31.61 – 3.10	Depositor EDS
% Data completeness (in resolution range)	93.0 (31.61-3.10) 92.9 (31.61-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 3.13Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.195 , 0.256 0.216 , 0.278	Depositor DCC
R_{free} test set	1008 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	46.9	Xtriage
Anisotropy	0.351	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 71.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	5883	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.43% 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: JLS, NAG

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	0.80	0/1288	1.16	2/1759 (0.1%)
2	B	0.74	0/1705	1.06	0/2352
3	C	0.74	0/2153	1.19	5/2949 (0.2%)
4	D	0.78	0/806	1.15	1/1101 (0.1%)
All	All	0.76	0/5952	1.14	8/8161 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	151	ASP	CA-CB-CG	7.71	120.31	112.60
1	A	7	ASP	CA-CB-CG	7.12	119.72	112.60
1	A	9	GLY	N-CA-C	5.38	118.74	112.23
3	C	86	LYS	CA-C-N	5.24	127.25	120.44
3	C	86	LYS	C-N-CA	5.24	127.25	120.44
3	C	75	SER	CA-C-N	5.22	127.22	120.44
3	C	75	SER	C-N-CA	5.22	127.22	120.44
4	D	29	GLY	N-CA-C	5.20	125.51	113.18

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	1265	0	1050	8	0
2	B	1658	0	1363	17	0
3	C	2088	0	1946	24	0
4	D	783	0	710	11	0
5	C	28	0	26	0	0
6	C	58	0	93	1	0
7	C	2	0	0	0	0
7	D	1	0	0	0	0
All	All	5883	0	5188	56	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 (56) 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)
4:D:39:LEU:HD23	4:D:80:CYS:HB3	1.41	0.98
3:C:88:LEU:HD23	3:C:143:LEU:HD23	1.66	0.78
4:D:39:LEU:CD2	4:D:80:CYS:HB3	2.18	0.73
2:B:21:LEU:CD1	2:B:41:TRP:CH2	2.75	0.70
1:A:143:ASP:HA	2:B:139:PHE:HA	1.76	0.67
4:D:39:LEU:HD23	4:D:80:CYS:CB	2.21	0.66
3:C:209:SER:HB3	3:C:244:TYR:HD1	1.61	0.65
3:C:98:VAL:HG22	3:C:116:VAL:HG22	1.83	0.60
3:C:187:PRO:HB3	3:C:211:PHE:HB3	1.87	0.56
2:B:107:SER:HB2	2:B:114:GLN:HG2	1.87	0.55
2:B:43:ARG:HH21	2:B:100:SER:HB2	1.72	0.55
3:C:133:PRO:HB3	3:C:145:ILE:HD11	1.89	0.53
3:C:218:VAL:HG22	3:C:263:VAL:HG13	1.91	0.52
3:C:15:ILE:HG12	4:D:62:PHE:HE1	1.75	0.52
3:C:217:TRP:HB3	3:C:264:LYS:HB2	1.90	0.52
3:C:88:LEU:HD22	3:C:140:TRP:HB2	1.92	0.52
1:A:40:MET:HE3	2:B:113:PRO:HA	1.92	0.51
4:D:49:VAL:HG12	4:D:68:THR:HB	1.93	0.51
3:C:9:PRO:HB3	3:C:103:GLU:HB3	1.92	0.51
1:A:45:TYR:HB2	1:A:48:LYS:HD2	1.93	0.50
3:C:98:VAL:HG13	3:C:116:VAL:HG22	1.95	0.49
4:D:25:CYS:HB3	4:D:39:LEU:HD21	1.94	0.48
3:C:211:PHE:HB2	3:C:265:HIS:CE1	2.49	0.48
1:A:6:GLN:HG2	1:A:121:ASP:H	1.80	0.46
4:D:1:ILE:HG13	4:D:86:THR:HG22	1.97	0.46
4:D:33:SER:HB2	4:D:54:LEU:HD21	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:203:LEU:HD11	3:C:248:THR:HB	1.97	0.46
1:A:131:ILE:HD12	1:A:158:ASP:HA	1.97	0.45
1:A:32:TYR:HD2	1:A:57:TYR:HB3	1.80	0.45
3:C:14:GLN:HB3	3:C:98:VAL:HB	1.98	0.45
3:C:77:PHE:O	3:C:81:VAL:HG23	2.17	0.45
2:B:55:TYR:CZ	2:B:67:ASP:HB2	2.52	0.45
2:B:53:ILE:HG22	2:B:54:HIS:ND1	2.31	0.44
3:C:212:TYR:CD1	3:C:213:PRO:HA	2.53	0.44
2:B:23:CYS:HB2	2:B:41:TRP:CZ2	2.53	0.43
4:D:25:CYS:CB	4:D:39:LEU:HD21	2.48	0.43
1:A:177:CYS:HB3	2:B:182:CYS:SG	2.58	0.43
2:B:112:GLN:HE21	2:B:112:GLN:N	2.16	0.43
3:C:148:LEU:HD11	6:C:303:JLS:H80	2.00	0.43
4:D:1:ILE:HD12	4:D:86:THR:HA	1.99	0.43
3:C:49:SER:HB3	3:C:54:SER:HB2	2.01	0.43
2:B:43:ARG:HB3	2:B:53:ILE:HD11	2.01	0.43
3:C:95:GLU:C	3:C:96:LEU:HD12	2.44	0.43
2:B:21:LEU:HD11	2:B:41:TRP:CZ3	2.55	0.42
3:C:205:VAL:HG22	3:C:248:THR:HG22	2.01	0.42
1:A:26:SER:HA	1:A:27:ASN:HA	1.75	0.42
2:B:21:LEU:HD11	2:B:41:TRP:CH2	2.54	0.42
3:C:88:LEU:CD2	3:C:143:LEU:HD23	2.41	0.42
2:B:99:THR:HG23	2:B:123:SER:HA	2.02	0.41
3:C:133:PRO:HB3	3:C:145:ILE:CD1	2.50	0.41
3:C:133:PRO:HD3	3:C:145:ILE:HD13	2.02	0.41
2:B:213:GLN:HG3	2:B:254:ALA:HA	2.02	0.41
4:D:37:VAL:HG13	4:D:82:VAL:HG22	2.02	0.41
2:B:102:TYR:CE2	2:B:122:LEU:HD22	2.56	0.41
2:B:138:VAL:HG23	2:B:248:ALA:HB3	2.02	0.41
3:C:106:PRO:HA	3:C:107:GLY:HA2	1.72	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	186/209 (89%)	170 (91%)	15 (8%)	1 (0%)	24	57
2	B	238/243 (98%)	226 (95%)	12 (5%)	0	100	100
3	C	268/278 (96%)	255 (95%)	13 (5%)	0	100	100
4	D	97/119 (82%)	93 (96%)	4 (4%)	0	100	100
All	All	789/849 (93%)	744 (94%)	44 (6%)	1 (0%)	48	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8	PRO

5.3.2 Protein sidechains ⓘ

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	114/188 (61%)	109 (96%)	5 (4%)	25	56
2	B	144/210 (69%)	142 (99%)	2 (1%)	59	76
3	C	217/243 (89%)	209 (96%)	8 (4%)	30	61
4	D	83/109 (76%)	80 (96%)	3 (4%)	31	62
All	All	558/750 (74%)	540 (97%)	18 (3%)	34	64

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASP
1	A	57	TYR
1	A	59	SER
1	A	110	LEU
1	A	179	LEU
2	B	25	GLN
2	B	112	GLN

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Mol	Chain	Res	Type
3	C	43	ASP
3	C	65	THR
3	C	67	GLN
3	C	123	ILE
3	C	124	LEU
3	C	165	THR
3	C	240	ASP
3	C	273	ILE
4	D	1	ILE
4	D	12	ARG
4	D	64	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	132	GLN
2	B	11	GLN
2	B	25	GLN
2	B	112	GLN
3	C	67	GLN
3	C	112	ASN
3	C	159	GLN
4	D	42	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

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
6	JLS	C	303	-	58,58,58	0.47	1 (1%)	64,67,67	0.71	0
5	NAG	C	302	3	14,14,15	0.30	0	17,19,21	0.74	1 (5%)
5	NAG	C	301	3	14,14,15	0.30	0	17,19,21	0.78	1 (5%)

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
6	JLS	C	303	-	-	22/56/76/76	0/1/1/1
5	NAG	C	302	3	-	0/6/23/26	0/1/1/1
5	NAG	C	301	3	-	1/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	303	JLS	O1A-C1A	2.14	1.43	1.40

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	301	NAG	C1-O5-C5	2.75	115.87	112.19
5	C	302	NAG	C1-O5-C5	2.32	115.29	112.19

There are no chirality outliers.

All (23) torsion outliers are listed below:

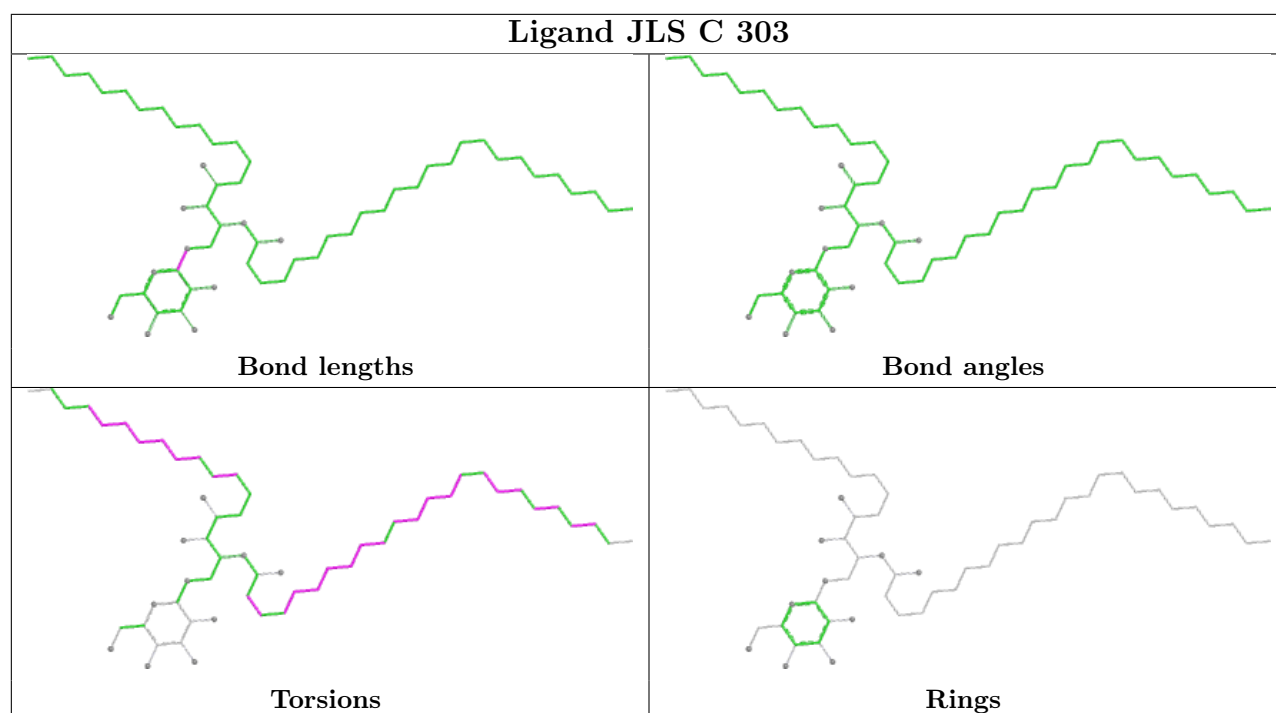
Mol	Chain	Res	Type	Atoms
6	C	303	JLS	CAA-CAB-CAC-CAD
6	C	303	JLS	CAJ-CAK-CAL-CAM
6	C	303	JLS	CAE-CAF-CAG-CAH
6	C	303	JLS	CAF-CAG-CAH-CAI
6	C	303	JLS	C9-C10-C11-C12
6	C	303	JLS	C11-C10-C9-C8
6	C	303	JLS	CAK-CAL-CAM-CAN
6	C	303	JLS	C13-C14-C15-C16
6	C	303	JLS	CAH-CAI-CAJ-CAK
6	C	303	JLS	C11-C12-C13-C14
6	C	303	JLS	CAD-CAE-CAF-CAG
6	C	303	JLS	CAP-CAQ-CAR-CAS
6	C	303	JLS	C6-C7-C8-C9
6	C	303	JLS	CAC-CAD-CAE-CAF
6	C	303	JLS	C12-C13-C14-C15
6	C	303	JLS	CAT-CAU-CAV-CAW
6	C	303	JLS	CAG-CAH-CAI-CAJ
6	C	303	JLS	C10-C11-C12-C13
6	C	303	JLS	CAR-CAS-CAT-CAU
5	C	301	NAG	O5-C5-C6-O6
6	C	303	JLS	CAL-CAM-CAN-CAO
6	C	303	JLS	CAM-CAN-CAO-CAP
6	C	303	JLS	CAO-CAP-CAQ-CAR

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	303	JLS	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	188/209 (89%)	0.78	25 (13%) 7 4	23, 74, 150, 159	0
2	B	240/243 (98%)	0.58	20 (8%) 17 9	21, 82, 153, 158	0
3	C	270/278 (97%)	-0.10	4 (1%) 72 52	16, 37, 96, 125	0
4	D	99/119 (83%)	0.11	3 (3%) 52 31	16, 53, 104, 112	0
All	All	797/849 (93%)	0.34	52 (6%) 25 13	16, 58, 144, 159	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	137	ALA	5.0
1	A	139	TYR	4.9
4	D	1	ILE	4.0
2	B	139	PHE	3.8
1	A	198	SER	3.6
2	B	143	GLU	3.5
1	A	136	PRO	3.3
1	A	200	PHE	3.2
2	B	196	ASP	3.2
1	A	5	GLU	3.0
1	A	143	ASP	2.9
1	A	158	ASP	2.9
1	A	168	ASP	2.9
1	A	162	ASN	2.9
2	B	208	SER	2.8
1	A	165	GLN	2.8
1	A	207	ASN	2.8
1	A	155	THR	2.8
3	C	202	LEU	2.7
2	B	140	GLU	2.6
2	B	230	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	138	VAL	2.6
2	B	141	PRO	2.6
2	B	153	THR	2.5
2	B	142	SER	2.5
1	A	6	GLN	2.5
3	C	107	GLY	2.5
2	B	131	VAL	2.4
1	A	133	ASN	2.4
1	A	170	ASP	2.4
1	A	167	LYS	2.4
4	D	99	MET	2.4
2	B	212	TRP	2.4
2	B	184	ASP	2.3
2	B	255	ASP	2.3
1	A	159	SER	2.3
2	B	20	THR	2.3
2	B	210	THR	2.3
2	B	151	LYS	2.3
1	A	169	SER	2.2
1	A	166	SER	2.2
2	B	48	MET	2.1
1	A	145	LYS	2.1
2	B	211	PHE	2.1
3	C	261	CYS	2.1
4	D	73	THR	2.1
1	A	153	LEU	2.1
1	A	156	ASP	2.1
2	B	236	GLN	2.1
1	A	196	ASN	2.0
2	B	233	GLU	2.0
3	C	269	GLU	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 ⓘ

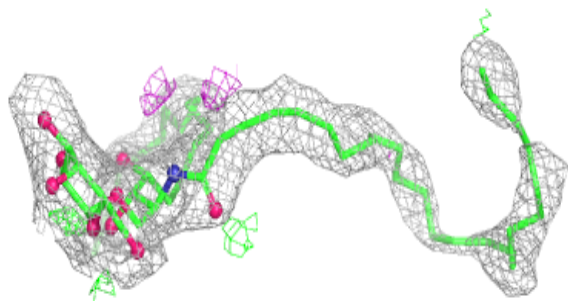
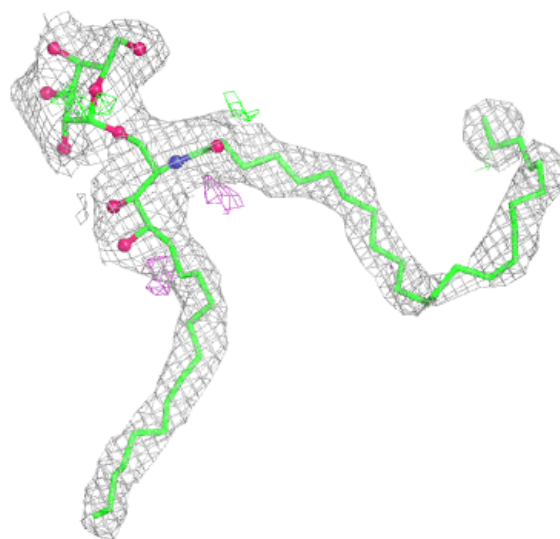
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	NAG	C	302	14/15	0.91	0.09	36,39,42,45	0
5	NAG	C	301	14/15	0.92	0.09	31,34,35,37	0
6	JLS	C	303	58/58	0.93	0.13	21,32,49,51	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around JLS C 303:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.