



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:48 AM UTC

PDB ID : 7RYN / pdb_00007ryn
Title : CD1a-sulfatide-gdTCR complex
Authors : Wegrecki, M.; Le Nours, J.; Rossjohn, J.
Deposited on : 2021-08-25
Resolution : 2.70 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 ⓘ](#)) 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

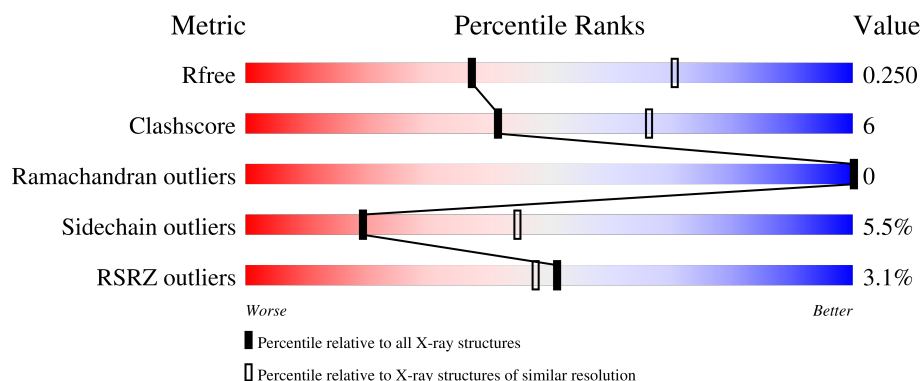
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	286	2% 81% 13% • 5%
2	B	108	2% 79% 13% 8%
3	C	248	4% 74% 19% • •
4	D	209	2% 67% 13% • 17%

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 6091 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 T-cell surface glycoprotein CD1a.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	0	0	0
			2164	1388	376	392	8			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ASP	-	expression tag	UNP P06126
A	0	ALA	-	expression tag	UNP P06126
A	2	THR	ASP	conflict	UNP P06126
A	13	ILE	THR	variant	UNP P06126
A	51	TRP	CYS	variant	UNP P06126
A	279	GLY	-	expression tag	UNP P06126
A	280	SER	-	expression tag	UNP P06126
A	281	LEU	-	expression tag	UNP P06126
A	282	VAL	-	expression tag	UNP P06126
A	283	PRO	-	expression tag	UNP P06126
A	284	ARG	-	expression tag	UNP P06126

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	0	0
			812	516	137	157	2			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	ASP	-	expression tag	UNP P61769
B	0	ALA	-	expression tag	UNP P61769
B	1	GLY	-	expression tag	UNP P61769
B	101	GLY	-	expression tag	UNP P61769
B	102	SER	-	expression tag	UNP P61769

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Chain	Residue	Modelled	Actual	Comment	Reference
B	103	LEU	-	expression tag	UNP P61769
B	104	VAL	-	expression tag	UNP P61769
B	105	PRO	-	expression tag	UNP P61769
B	106	ARG	-	expression tag	UNP P61769

- Molecule 3 is a protein called T cell receptor gamma variable 4,T cell receptor beta constant 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	238	Total	C	N	O	S	0	0	0
			1783	1134	296	347	6			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	MET	-	initiating methionine	UNP A0A0C4DH28
C	2	ALA	-	expression tag	UNP A0A0C4DH28
C	103	ASP	-	linker	UNP A0A0C4DH28
C	104	TYR	-	linker	UNP A0A0C4DH28
C	105	TYR	-	linker	UNP A0A0C4DH28
C	106	LYS	-	linker	UNP A0A0C4DH28
C	107	LYS	-	linker	UNP A0A0C4DH28
C	108	LEU	-	linker	UNP A0A0C4DH28
C	109	PHE	-	linker	UNP A0A0C4DH28
C	110	GLY	-	linker	UNP A0A0C4DH28
C	111	SER	-	linker	UNP A0A0C4DH28
C	112	GLY	-	linker	UNP A0A0C4DH28
C	113	THR	-	linker	UNP A0A0C4DH28
C	114	THR	-	linker	UNP A0A0C4DH28
C	115	LEU	-	linker	UNP A0A0C4DH28
C	116	VAL	-	linker	UNP A0A0C4DH28
C	117	VAL	-	linker	UNP A0A0C4DH28
C	118	THR	-	linker	UNP A0A0C4DH28
C	119	GLU	-	linker	UNP A0A0C4DH28
C	122	LYS	ASN	conflict	UNP P01850
C	123	ASN	LYS	conflict	UNP P01850
C	155	TYR	PHE	conflict	UNP P01850
C	175	CYS	SER	engineered mutation	UNP P01850
C	193	ALA	CYS	engineered mutation	UNP P01850

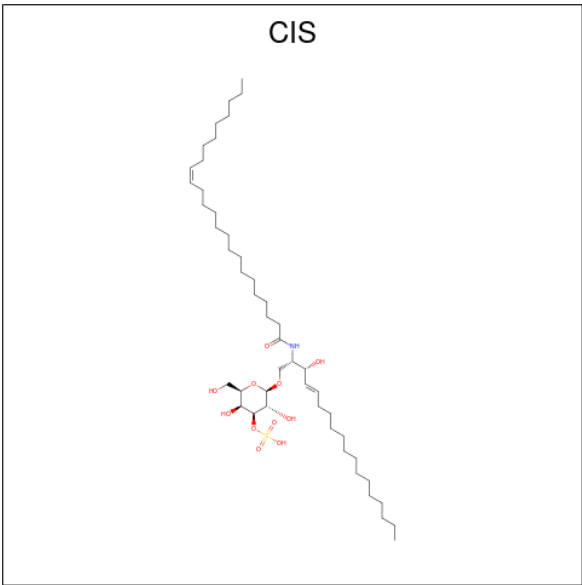
- Molecule 4 is a protein called T cell receptor delta variable 1,T cell receptor alpha chain constant.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	173	Total	C	N	O	S	0	0	0
			1202	763	201	233	5			

There are 22 discrepancies between the modelled and reference sequences:

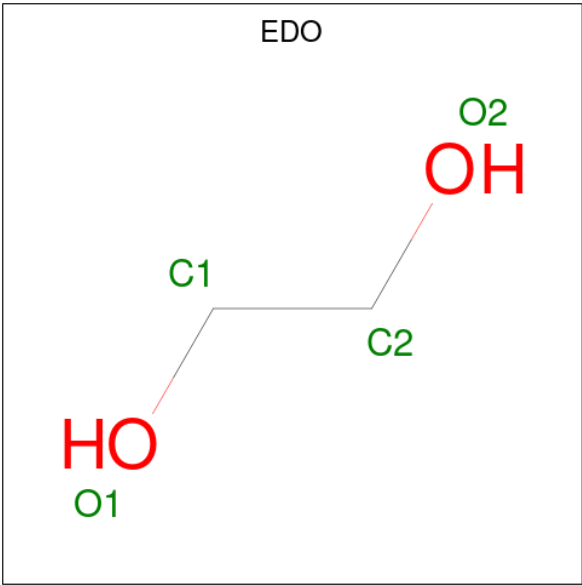
Chain	Residue	Modelled	Actual	Comment	Reference
D	1	MET	-	initiating methionine	UNP A0A1B0GX56
D	97	LEU	-	linker	UNP A0A1B0GX56
D	98	ARG	-	linker	UNP A0A1B0GX56
D	99	TRP	-	linker	UNP A0A1B0GX56
D	100	PRO	-	linker	UNP A0A1B0GX56
D	101	ASP	-	linker	UNP A0A1B0GX56
D	102	LYS	-	linker	UNP A0A1B0GX56
D	103	LEU	-	linker	UNP A0A1B0GX56
D	104	ILE	-	linker	UNP A0A1B0GX56
D	105	PHE	-	linker	UNP A0A1B0GX56
D	106	GLY	-	linker	UNP A0A1B0GX56
D	107	LYS	-	linker	UNP A0A1B0GX56
D	108	GLY	-	linker	UNP A0A1B0GX56
D	109	THR	-	linker	UNP A0A1B0GX56
D	110	ARG	-	linker	UNP A0A1B0GX56
D	111	VAL	-	linker	UNP A0A1B0GX56
D	112	THR	-	linker	UNP A0A1B0GX56
D	113	VAL	-	linker	UNP A0A1B0GX56
D	114	GLU	-	linker	UNP A0A1B0GX56
D	115	PRO	-	linker	UNP A0A1B0GX56
D	116	ASN	-	linker	UNP A0A1B0GX56
D	163	CYS	THR	engineered mutation	UNP P01848

- Molecule 5 is (15Z)-N-((1S,2R,3E)-2-HYDROXY-1-[(3-O-SULFO-BETA-D-GALACTO PYRANOSYL)OXY]METHYL}HEPTADEC-3-ENYL)TETRACOS-15-ENAMIDE (CCD ID: CIS) (formula: C₄₈H₉₁NO₁₁S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			61	48	1	11	1		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



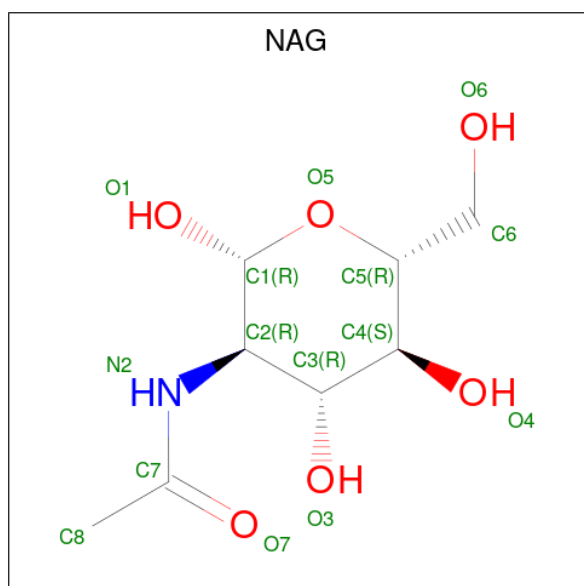
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		

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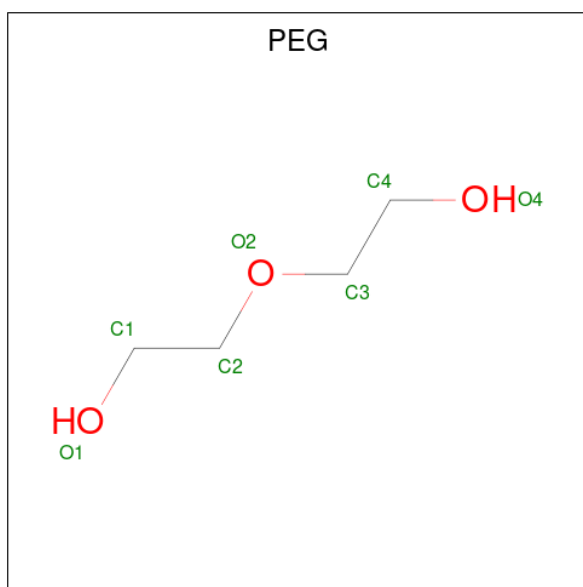
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



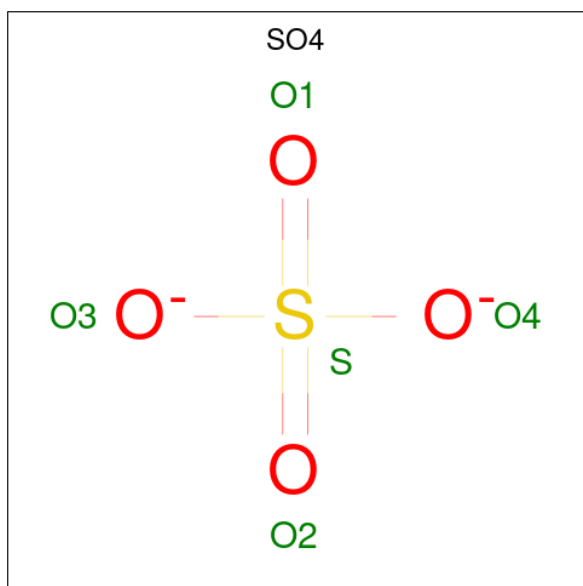
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 8 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 9 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	O	S	0	0
			5	4	1		
9	C	1	Total	O	S	0	0
			5	4	1		

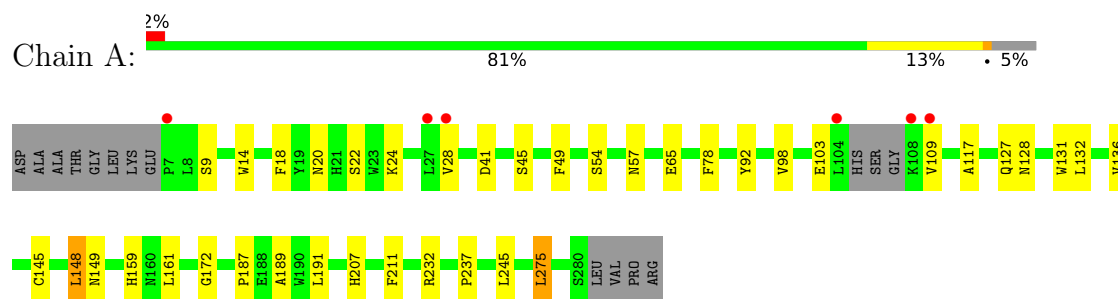
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	8	Total 8	O 8	0	0
10	B	6	Total 6	O 6	0	0

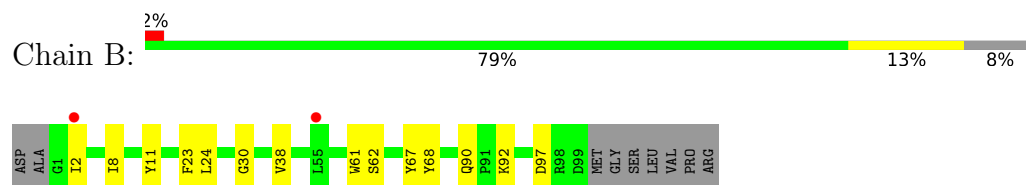
3 Residue-property plots [i](#)

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.

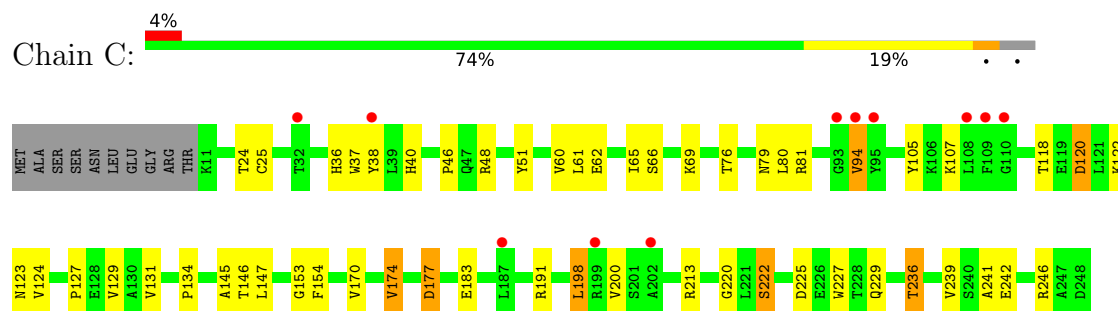
- Molecule 1: T-cell surface glycoprotein CD1a



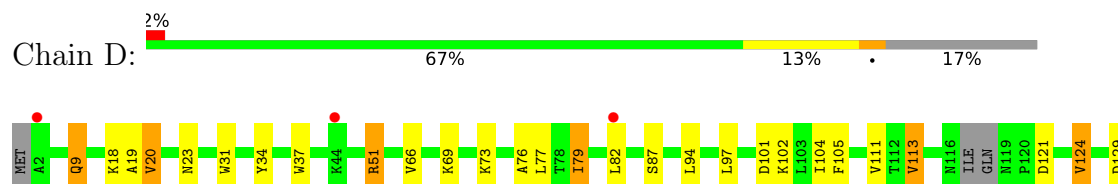
- Molecule 2: Beta-2-microglobulin

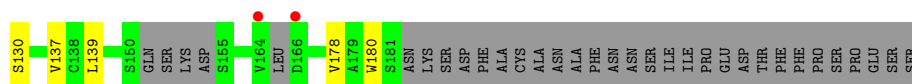


- Molecule 3: T cell receptor gamma variable 4,T cell receptor beta constant 1



- Molecule 4: T cell receptor delta variable 1,T cell receptor alpha chain constant





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	234.63Å 42.21Å 123.06Å 90.00° 114.29° 90.00°	Depositor
Resolution (Å)	42.93 – 2.70 42.93 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (42.93-2.70) 99.9 (42.93-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.80 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.232 , 0.268 (Not available) , 0.250	Depositor DCC
R_{free} test set	1569 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	60.0	Xtriage
Anisotropy	0.767	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 66.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6091	wwPDB-VP
Average B, all atoms (Å ²)	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.34% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, NAG, CIS, SO4, PEG

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	0.13	0/2235	0.33	0/3047
2	B	0.11	0/835	0.29	0/1135
3	C	0.12	0/1831	0.32	0/2512
4	D	0.13	0/1226	0.31	0/1682
All	All	0.13	0/6127	0.32	0/8376

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	A	2164	0	2015	24	0
2	B	812	0	755	7	0
3	C	1783	0	1618	27	0
4	D	1202	0	1004	21	0
5	A	61	0	91	3	0
6	A	20	0	30	0	0
6	B	4	0	6	1	0
7	A	14	0	13	0	0
8	A	7	0	10	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	A	5	0	0	0	0
9	C	5	0	0	0	0
10	A	8	0	0	0	0
10	B	6	0	0	0	0
All	All	6091	0	5542	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 6.

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:187:PRO:HB3	1:A:211:PHE:HB3	1.65	0.78
1:A:14:TRP:HD1	1:A:28:VAL:HG22	1.53	0.74
4:D:20:VAL:HG12	4:D:79:ILE:HD11	1.71	0.72
3:C:213:ARG:HG3	3:C:242:GLU:HB3	1.76	0.68
1:A:161:LEU:HD23	5:A:301:CIS:H281	1.77	0.67
4:D:69:LYS:O	4:D:73:LYS:N	2.28	0.67
3:C:40:HIS:HB3	3:C:94:VAL:HG13	1.78	0.65
4:D:124:VAL:HA	4:D:139:LEU:O	1.97	0.63
3:C:153:GLY:HA2	3:C:191:ARG:HD3	1.80	0.63
3:C:134:PRO:HG2	3:C:145:ALA:HB1	1.80	0.63
4:D:9:GLN:HE22	4:D:23:ASN:H	1.45	0.62
3:C:120:ASP:HB3	3:C:122:LYS:HG2	1.82	0.61
1:A:145:CYS:O	1:A:149:ASN:ND2	2.34	0.61
2:B:8:ILE:HD12	2:B:92:LYS:HD2	1.82	0.60
3:C:174:VAL:HB	3:C:198:LEU:HD23	1.85	0.58
3:C:66:SER:HB2	3:C:69:LYS:HB2	1.86	0.58
3:C:38:TYR:OH	3:C:107:LYS:NZ	2.38	0.57
1:A:49:PHE:HB3	1:A:54:SER:HB2	1.88	0.55
1:A:237:PRO:O	2:B:11:TYR:OH	2.22	0.55
3:C:105:TYR:HA	4:D:51:ARG:HH21	1.72	0.55
3:C:36:HIS:HD2	3:C:48:ARG:HH11	1.54	0.55
4:D:66:VAL:HA	4:D:76:ALA:O	2.07	0.54
3:C:220:GLY:H	3:C:236:THR:HB	1.73	0.54
4:D:79:ILE:HD13	4:D:82:LEU:HD13	1.91	0.52
3:C:62:GLU:HB3	3:C:65:ILE:HD13	1.91	0.52
3:C:134:PRO:HD3	3:C:147:LEU:HD13	1.90	0.52
1:A:148:LEU:HD11	5:A:301:CIS:H82	1.92	0.52
3:C:24:THR:HG22	3:C:81:ARG:HG3	1.92	0.51
2:B:30:GLY:HA2	2:B:62:SER:HB2	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:79:ASN:HB3	3:C:81:ARG:NH1	2.27	0.50
4:D:34:TYR:CZ	4:D:97:LEU:HG	2.48	0.49
4:D:129:ASP:OD1	4:D:130:SER:N	2.46	0.48
1:A:41:ASP:O	1:A:45:SER:HA	2.13	0.48
1:A:22:SER:OG	4:D:101:ASP:OD1	2.26	0.48
1:A:128:ASN:O	1:A:159:HIS:ND1	2.45	0.48
3:C:46:PRO:HG2	4:D:105:PHE:CG	2.49	0.47
4:D:18:LYS:HG3	4:D:19:ALA:H	1.79	0.47
2:B:23:PHE:CE2	6:B:401:EDO:H11	2.50	0.47
4:D:87:SER:HA	4:D:111:VAL:HG23	1.97	0.47
1:A:127:GLN:HG3	1:A:132:LEU:HD22	1.96	0.47
3:C:134:PRO:HB3	3:C:146:THR:H	1.80	0.46
1:A:41:ASP:O	1:A:45:SER:N	2.48	0.46
3:C:76:THR:HG22	3:C:80:LEU:HD13	1.97	0.46
3:C:25:CYS:HB3	3:C:37:TRP:CH2	2.51	0.46
4:D:87:SER:HB2	4:D:113:VAL:HG12	1.98	0.45
4:D:37:TRP:CE2	4:D:77:LEU:HB2	2.52	0.45
4:D:31:TRP:CD1	4:D:31:TRP:H	2.33	0.45
1:A:41:ASP:O	1:A:45:SER:CA	2.65	0.44
4:D:101:ASP:CG	4:D:102:LYS:H	2.26	0.44
4:D:121:ASP:N	4:D:121:ASP:OD1	2.50	0.44
1:A:18:PHE:HB3	1:A:92:TYR:CE2	2.53	0.44
1:A:131:TRP:HZ2	5:A:301:CIS:H232	1.82	0.44
1:A:232:ARG:HG3	1:A:245:LEU:HD11	1.99	0.44
3:C:127:PRO:HB3	3:C:154:PHE:HB3	1.99	0.44
3:C:131:VAL:HG23	3:C:241:ALA:HB3	2.00	0.44
4:D:94:LEU:HD23	4:D:104:ILE:HB	2.00	0.43
1:A:57:ASN:HB3	1:A:172:GLY:HA3	2.01	0.43
1:A:14:TRP:CD1	1:A:28:VAL:HG22	2.43	0.43
2:B:38:VAL:HB	2:B:67:TYR:CZ	2.54	0.43
3:C:177:ASP:OD1	3:C:177:ASP:N	2.39	0.42
1:A:131:TRP:CD2	1:A:148:LEU:HD23	2.54	0.42
3:C:222:SER:O	3:C:225:ASP:HB2	2.19	0.42
3:C:129:VAL:HG12	3:C:239:VAL:HG12	2.00	0.42
4:D:137:VAL:HG12	4:D:180:TRP:HB3	2.01	0.42
4:D:9:GLN:NE2	4:D:23:ASN:H	2.15	0.41
1:A:191:LEU:HD11	1:A:275:LEU:HB3	2.02	0.41
3:C:51:TYR:HB3	3:C:60:VAL:HG13	2.01	0.41
3:C:134:PRO:HG3	3:C:146:THR:C	2.45	0.41
1:A:14:TRP:HB3	1:A:98:VAL:HB	2.01	0.41
3:C:227:TRP:CZ2	3:C:229:GLN:HB2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:SER:HB3	1:A:103:GLU:HG2	2.02	0.41
1:A:189:ALA:HA	1:A:207:HIS:O	2.20	0.41
2:B:24:LEU:O	2:B:68:TYR:HA	2.21	0.40
1:A:117:ALA:HB2	2:B:61:TRP:CE2	2.56	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	267/286 (93%)	259 (97%)	8 (3%)	0	100	100
2	B	97/108 (90%)	95 (98%)	2 (2%)	0	100	100
3	C	236/248 (95%)	227 (96%)	9 (4%)	0	100	100
4	D	165/209 (79%)	154 (93%)	11 (7%)	0	100	100
All	All	765/851 (90%)	735 (96%)	30 (4%)	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	A	228/247 (92%)	220 (96%)	8 (4%)	32	61
2	B	90/100 (90%)	87 (97%)	3 (3%)	33	63

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	180/215 (84%)	165 (92%)	15 (8%)	10	26
4	D	104/188 (55%)	97 (93%)	7 (7%)	15	36
All	All	602/750 (80%)	569 (94%)	33 (6%)	19	45

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	ASN
1	A	24	LYS
1	A	65	GLU
1	A	78	PHE
1	A	109	VAL
1	A	136	VAL
1	A	148	LEU
1	A	275	LEU
2	B	2	ILE
2	B	90	GLN
2	B	97	ASP
3	C	61	LEU
3	C	94	VAL
3	C	118	THR
3	C	120	ASP
3	C	123	ASN
3	C	124	VAL
3	C	170	VAL
3	C	174	VAL
3	C	177	ASP
3	C	183	GLU
3	C	198	LEU
3	C	200	VAL
3	C	222	SER
3	C	236	THR
3	C	246	ARG
4	D	9	GLN
4	D	20	VAL
4	D	51	ARG
4	D	79	ILE
4	D	113	VAL
4	D	124	VAL
4	D	178	VAL

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

Mol	Chain	Res	Type
1	A	129	ASN
2	B	18	ASN
3	C	16	GLN
3	C	36	HIS
3	C	79	ASN
3	C	86	ASN
3	C	90	ASN
3	C	123	ASN
3	C	215	GLN
3	C	217	GLN

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 ⓘ

11 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	EDO	A	302	-	3,3,3	0.43	0	2,2,2	0.37	0
6	EDO	A	304	-	3,3,3	0.44	0	2,2,2	0.33	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	SO4	C	301	-	4,4,4	0.22	0	6,6,6	0.08	0
8	PEG	A	308	-	6,6,6	0.10	0	5,5,5	0.14	0
7	NAG	A	303	1	14,14,15	0.25	0	17,19,21	0.52	0
6	EDO	B	401	-	3,3,3	0.42	0	2,2,2	0.38	0
6	EDO	A	305	-	3,3,3	0.43	0	2,2,2	0.32	0
6	EDO	A	306	-	3,3,3	0.44	0	2,2,2	0.32	0
9	SO4	A	309	-	4,4,4	0.30	0	6,6,6	0.10	0
6	EDO	A	307	-	3,3,3	0.42	0	2,2,2	0.41	0
5	CIS	A	301	-	60,61,61	0.24	0	68,72,72	0.59	2 (2%)

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	EDO	A	302	-	-	0/1/1/1	-
6	EDO	A	304	-	-	0/1/1/1	-
8	PEG	A	308	-	-	2/4/4/4	-
7	NAG	A	303	1	-	0/6/23/26	0/1/1/1
6	EDO	B	401	-	-	1/1/1/1	-
6	EDO	A	305	-	-	0/1/1/1	-
6	EDO	A	306	-	-	0/1/1/1	-
6	EDO	A	307	-	-	0/1/1/1	-
5	CIS	A	301	-	-	22/58/78/78	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	301	CIS	O8-S-O9	3.36	125.20	112.24
5	A	301	CIS	O10-S-O4	-3.12	99.21	106.37

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	301	CIS	C-C1-C2-C3
5	A	301	CIS	N-C1-C2-C3
5	A	301	CIS	C-C1-C2-O1

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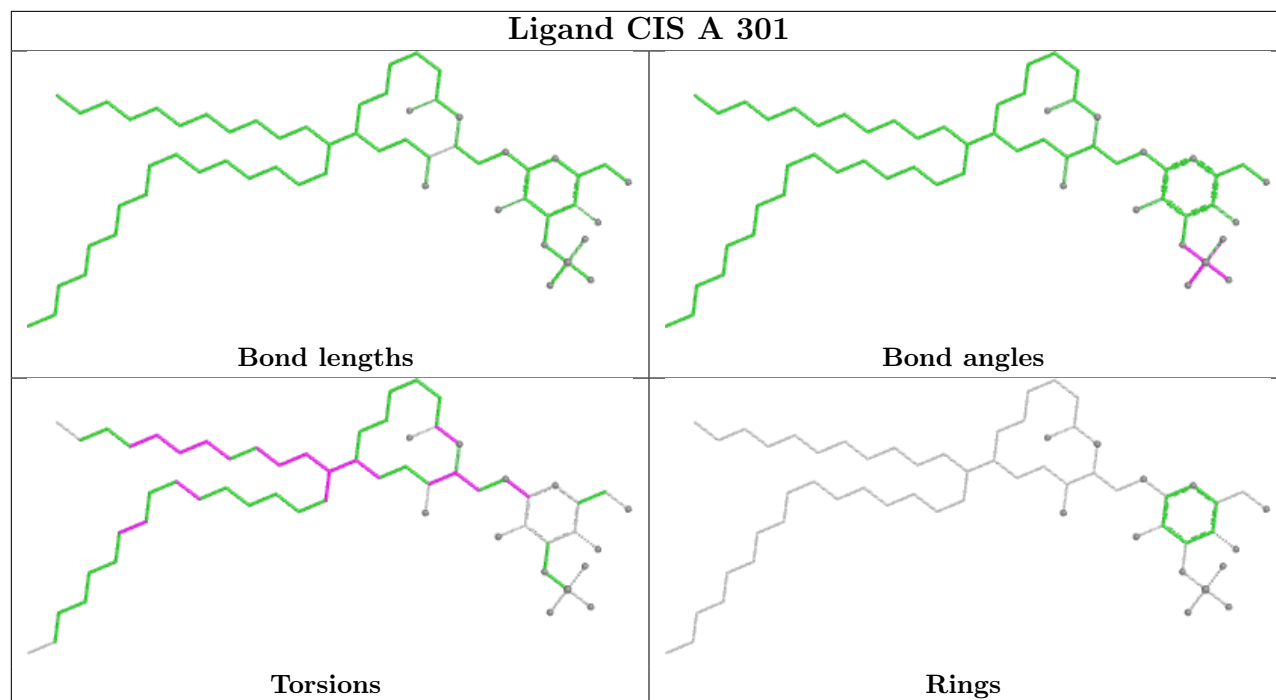
Mol	Chain	Res	Type	Atoms
5	A	301	CIS	N-C1-C2-O1
5	A	301	CIS	O-C-C1-C2
5	A	301	CIS	O-C-C1-N
5	A	301	CIS	O6-C43-O-C
5	A	301	CIS	C19-C18-N-C1
5	A	301	CIS	C44-C43-O-C
5	A	301	CIS	O2-C18-N-C1
5	A	301	CIS	C10-C11-C12-C13
5	A	301	CIS	C23-C24-C25-C26
5	A	301	CIS	C5-C6-C7-C8
5	A	301	CIS	C9-C10-C11-C12
5	A	301	CIS	C24-C25-C26-C27
5	A	301	CIS	C3-C4-C5-C6
5	A	301	CIS	C11-C12-C13-C14
5	A	301	CIS	C12-C13-C14-C15
8	A	308	PEG	O1-C1-C2-O2
8	A	308	PEG	O2-C3-C4-O4
5	A	301	CIS	C7-C8-C9-C10
5	A	301	CIS	C6-C7-C8-C9
5	A	301	CIS	C30-C31-C32-C33
6	B	401	EDO	O1-C1-C2-O2
5	A	301	CIS	C33-C34-C35-C36

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	401	EDO	1	0
5	A	301	CIS	3	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	271/286 (94%)	-0.03	6 (2%) 62 59	47, 71, 101, 125	0
2	B	99/108 (91%)	-0.09	2 (2%) 65 62	44, 59, 93, 103	0
3	C	238/248 (95%)	0.44	11 (4%) 37 34	71, 126, 176, 209	0
4	D	173/209 (82%)	0.50	5 (2%) 53 50	72, 142, 202, 226	0
All	All	781/851 (91%)	0.22	24 (3%) 51 48	44, 91, 177, 226	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	94	VAL	4.2
1	A	109	VAL	4.0
3	C	108	LEU	3.7
1	A	108	LYS	3.6
3	C	109	PHE	3.4
3	C	93	GLY	3.1
1	A	7	PRO	3.0
1	A	104	LEU	3.0
1	A	28	VAL	2.9
4	D	164	VAL	2.7
3	C	199	ARG	2.6
4	D	44	LYS	2.6
3	C	202	ALA	2.6
3	C	110	GLY	2.5
3	C	32	THR	2.5
2	B	55	LEU	2.5
4	D	166	ASP	2.4
4	D	82	LEU	2.3
3	C	95	TYR	2.3
4	D	2	ALA	2.2
3	C	38	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
3	C	187	LEU	2.1
2	B	2	ILE	2.1
1	A	27	LEU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

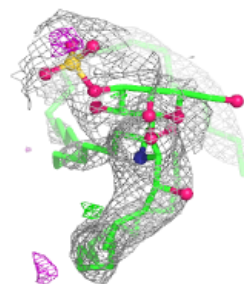
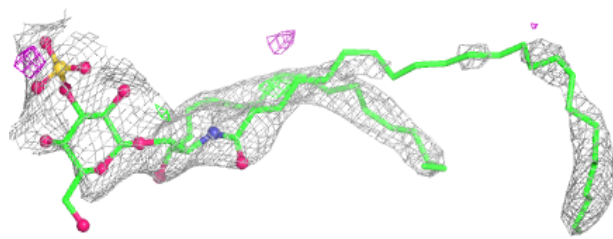
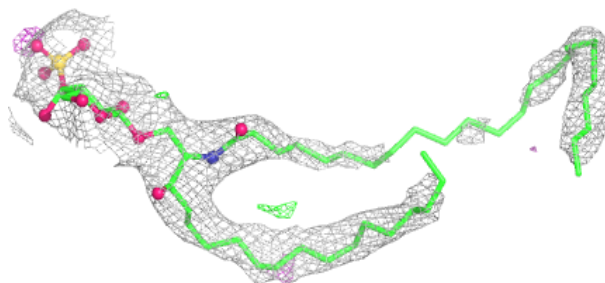
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	NAG	A	303	14/15	0.19	0.13	110,139,152,153	0
6	EDO	A	306	4/4	0.59	0.23	77,77,92,98	0
6	EDO	B	401	4/4	0.71	0.22	53,71,74,79	0
9	SO4	A	309	5/5	0.71	0.11	97,109,136,230	0
6	EDO	A	304	4/4	0.72	0.20	86,91,93,95	0
9	SO4	C	301	5/5	0.78	0.09	148,152,154,154	5
6	EDO	A	302	4/4	0.83	0.14	45,71,71,84	0
8	PEG	A	308	7/7	0.83	0.12	65,70,73,77	0
6	EDO	A	305	4/4	0.87	0.15	46,49,55,62	0
5	CIS	A	301	61/61	0.90	0.15	43,75,120,133	0
6	EDO	A	307	4/4	0.91	0.15	50,51,53,54	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 CIS A 301:

$2mF_o-DF_c$ (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.