



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 10:00 AM UTC

PDB ID : 7ZT4 / pdb_00007zt4
Title : Structure of E8 TCR in complex with human MR1 bound to 6FP
Authors : Karuppiah, V.; Robinson, R.A.
Deposited on : 2022-05-09
Resolution : 2.02 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

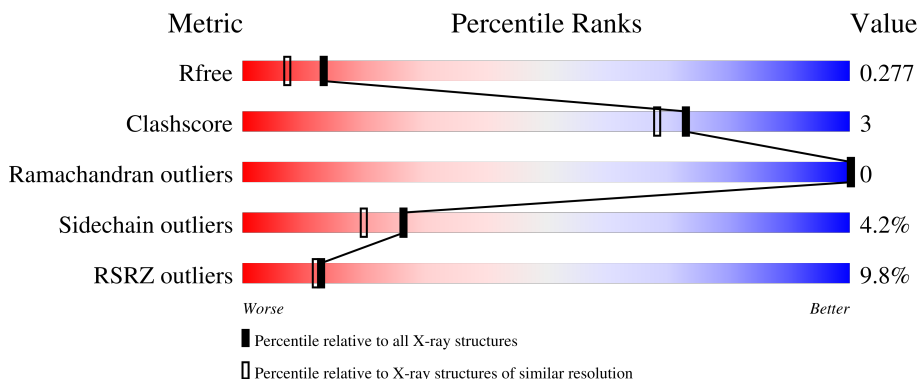
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.02 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	13299 (2.04-2.00)
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)
RSRZ outliers	180081	13314 (2.04-2.00)

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 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	290	12% 83% 8% 9%
2	B	100	7% 81% 17% ..
3	D	205	11% 74% 15% 11%
4	E	262	5% 83% 9% 8%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6637 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 Major histocompatibility complex class I-related gene protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	264	2192	1402	377	402	11	0	1	0

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q95460
A	261	SER	CYS	conflict	UNP Q95460
A	271	GLY	-	expression tag	UNP Q95460
A	272	SER	-	expression tag	UNP Q95460
A	273	GLY	-	expression tag	UNP Q95460
A	274	GLY	-	expression tag	UNP Q95460
A	275	GLY	-	expression tag	UNP Q95460
A	276	LEU	-	expression tag	UNP Q95460
A	277	ASN	-	expression tag	UNP Q95460
A	278	ASP	-	expression tag	UNP Q95460
A	279	ILE	-	expression tag	UNP Q95460
A	280	PHE	-	expression tag	UNP Q95460
A	281	GLU	-	expression tag	UNP Q95460
A	282	ALA	-	expression tag	UNP Q95460
A	283	GLN	-	expression tag	UNP Q95460
A	284	LYS	-	expression tag	UNP Q95460
A	285	ILE	-	expression tag	UNP Q95460
A	286	GLU	-	expression tag	UNP Q95460
A	287	TRP	-	expression tag	UNP Q95460
A	288	HIS	-	expression tag	UNP Q95460
A	289	GLU	-	expression tag	UNP Q95460

- 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
			828	528	140	157	3			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P61769

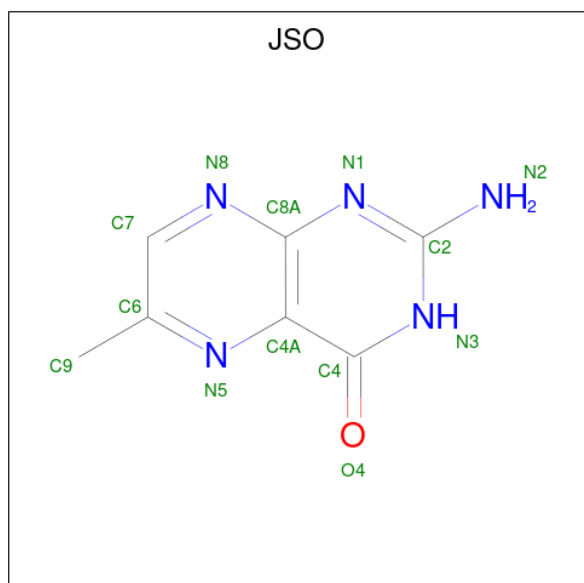
- Molecule 3 is a protein called TCR alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	183	Total	C	N	O	S	0	0	0
			1435	906	233	287	9			

- Molecule 4 is a protein called TCR beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	242	Total	C	N	O	S	0	2	0
			1935	1214	335	377	9			

- Molecule 5 is 2-azanyl-6-methyl-3 {H}-pteridin-4-one (CCD ID: JSO) (formula: C₇H₇N₅O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			13	7	5	1		

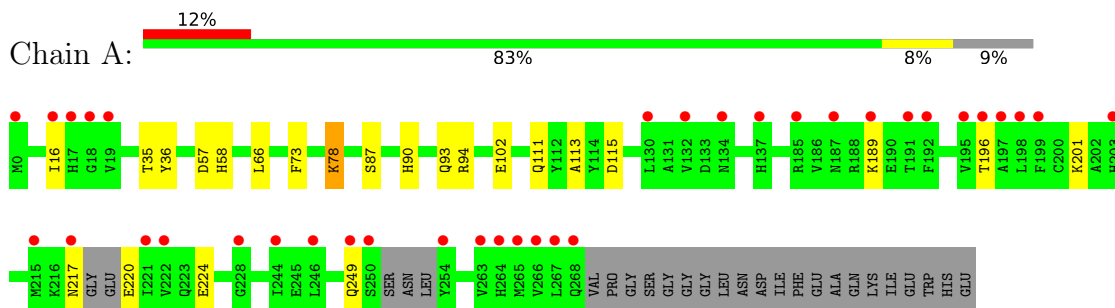
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	68	Total 68	O 68	0	0
6	B	27	Total 27	O 27	0	0
6	D	51	Total 51	O 51	0	0
6	E	88	Total 88	O 88	0	0

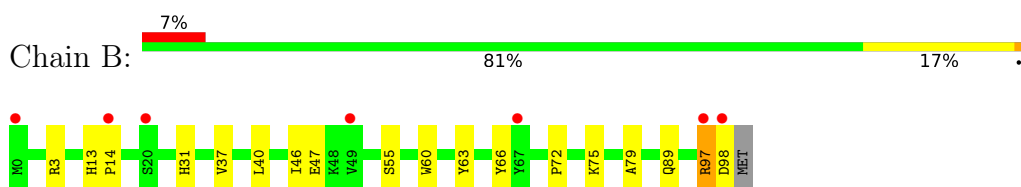
3 Residue-property plots [i](#)

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.

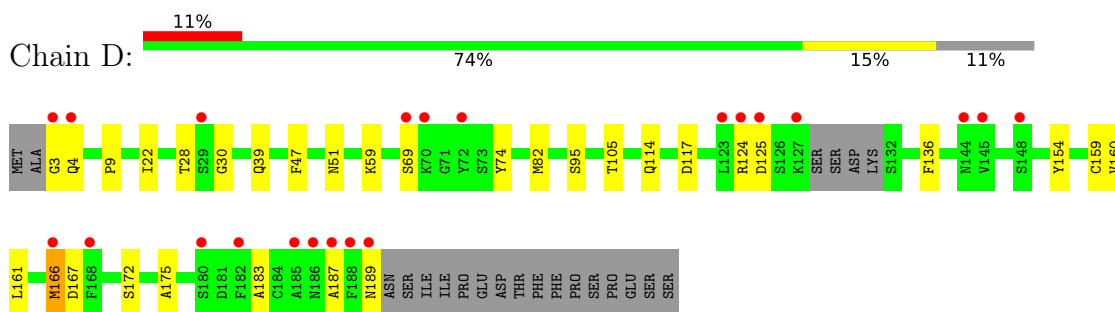
- Molecule 1: Major histocompatibility complex class I-related gene protein



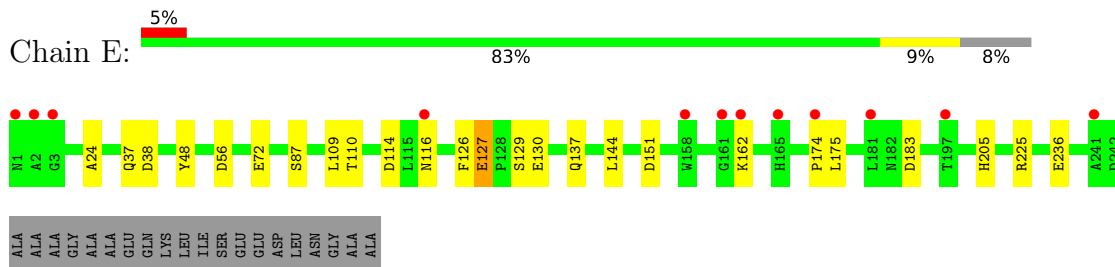
- Molecule 2: Beta-2-microglobulin



- Molecule 3: TCR alpha



- Molecule 4: TCR beta



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	91.71Å 104.39Å 117.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	58.99 – 2.02 58.99 – 2.02	Depositor EDS
% Data completeness (in resolution range)	100.0 (58.99-2.02) 99.9 (58.99-2.02)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.15 (at 2.02Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.234 , 0.268 0.244 , 0.277	Depositor DCC
R_{free} test set	3763 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	47.6	Xtriage
Anisotropy	0.390	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.1	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	0.95	EDS
Total number of atoms	6637	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.90% 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: JSO

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	1.06	1/2258 (0.0%)	1.31	3/3066 (0.1%)
2	B	1.03	0/851	1.20	0/1152
3	D	1.07	0/1465	1.26	2/1981 (0.1%)
4	E	1.07	3/1994 (0.2%)	1.28	3/2718 (0.1%)
All	All	1.06	4/6568 (0.1%)	1.28	8/8917 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	127	GLU	C-O	5.15	1.30	1.24
1	A	66	LEU	C-O	5.04	1.30	1.24
4	E	137[A]	GLN	C-O	5.03	1.32	1.24
4	E	137[B]	GLN	C-O	5.03	1.32	1.24

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	137[A]	GLN	CA-C-O	6.98	128.56	120.32
4	E	137[B]	GLN	CA-C-O	6.98	128.56	120.32
3	D	47	PHE	CA-C-O	-6.20	113.85	120.92
1	A	35	THR	CA-CB-OG1	-6.06	100.50	109.60
3	D	117	ASP	CB-CA-C	5.46	117.92	111.15
1	A	78	LYS	CA-C-N	5.29	127.68	120.54
1	A	78	LYS	C-N-CA	5.29	127.68	120.54
4	E	114	ASP	CA-CB-CG	5.07	117.67	112.60

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	2192	0	2077	8	0
2	B	828	0	794	10	0
3	D	1435	0	1358	15	0
4	E	1935	0	1827	14	0
5	A	13	0	0	0	0
6	A	68	0	0	0	0
6	B	27	0	0	1	0
6	D	51	0	0	0	0
6	E	88	0	0	0	0
All	All	6637	0	6056	40	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 3.

All (40) 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)
3:D:39:GLN:HE22	4:E:37:GLN:HE22	1.16	0.94
3:D:3:GLY:N	3:D:28:THR:HG1	1.81	0.78
4:E:151[B]:ASP:OD1	4:E:174:PRO:HG2	1.93	0.69
2:B:3:ARG:HD2	6:B:118:HOH:O	1.99	0.62
3:D:82:MET:HE2	3:D:160:VAL:HG11	1.83	0.61
1:A:102:GLU:H	1:A:102:GLU:CD	2.10	0.59
3:D:125:ASP:HA	4:E:126:PHE:CD2	2.39	0.58
4:E:205:HIS:HE1	4:E:236:GLU:OE1	1.88	0.57
2:B:72:PRO:HB2	2:B:97:ARG:HH12	1.72	0.55
3:D:166:MET:O	3:D:167:ASP:C	2.50	0.53
4:E:151[B]:ASP:CG	4:E:174:PRO:HG2	2.34	0.53
3:D:124:ARG:HB2	4:E:127:GLU:HB2	1.91	0.52
3:D:125:ASP:HA	4:E:126:PHE:HD2	1.73	0.52
1:A:90:HIS:HD2	1:A:115:ASP:OD2	1.95	0.50
3:D:30:GLY:HA3	3:D:95:SER:OG	2.13	0.49
4:E:48:TYR:CE1	4:E:56:ASP:HB3	2.48	0.48
1:A:87:SER:O	1:A:90:HIS:HE1	1.98	0.46
2:B:40:LEU:HB2	2:B:79:ALA:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:46:ILE:HG22	2:B:47:GLU:O	2.16	0.46
3:D:136:PHE:O	3:D:172:SER:HA	2.16	0.46
2:B:55:SER:HB3	2:B:63:TYR:CZ	2.52	0.45
3:D:9:PRO:O	3:D:105:THR:OG1	2.33	0.45
1:A:113:ALA:HB2	2:B:60:TRP:CE2	2.52	0.45
2:B:97:ARG:O	2:B:98:ASP:C	2.60	0.45
1:A:57:ASP:OD1	1:A:58:HIS:N	2.50	0.44
1:A:93:GLN:OE1	2:B:31:HIS:HE1	2.00	0.43
1:A:36:TYR:CD1	1:A:36:TYR:C	2.97	0.43
1:A:102:GLU:CD	1:A:102:GLU:N	2.76	0.42
2:B:37:VAL:HB	2:B:66:TYR:CZ	2.54	0.42
4:E:116:ASN:O	4:E:225:ARG:NH2	2.53	0.41
4:E:109:LEU:HD23	4:E:109:LEU:C	2.45	0.41
4:E:109:LEU:HD23	4:E:110:THR:N	2.35	0.41
2:B:13:HIS:O	2:B:14:PRO:C	2.63	0.41
3:D:3:GLY:N	3:D:28:THR:OG1	2.50	0.41
4:E:24:ALA:HA	4:E:72:GLU:O	2.21	0.41
3:D:183:ALA:O	3:D:187:ALA:HB2	2.21	0.41
3:D:154:TYR:HB3	4:E:175:LEU:HD23	2.03	0.40
3:D:22:ILE:O	3:D:74:TYR:HA	2.21	0.40
3:D:154:TYR:O	3:D:175:ALA:HA	2.22	0.40
4:E:38:ASP:OD1	4:E:87:SER:OG	2.38	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	259/290 (89%)	251 (97%)	8 (3%)	0	100	100
2	B	97/100 (97%)	91 (94%)	6 (6%)	0	100	100
3	D	179/205 (87%)	168 (94%)	11 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	E	242/262 (92%)	233 (96%)	9 (4%)	0	100	100
All	All	777/857 (91%)	743 (96%)	34 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	235/255 (92%)	223 (95%)	12 (5%)	21	14
2	B	94/95 (99%)	91 (97%)	3 (3%)	34	30
3	D	159/181 (88%)	150 (94%)	9 (6%)	18	11
4	E	213/222 (96%)	208 (98%)	5 (2%)	44	43
All	All	701/753 (93%)	672 (96%)	29 (4%)	26	20

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	ILE
1	A	73	PHE
1	A	78	LYS
1	A	94	ARG
1	A	111	GLN
1	A	189	LYS
1	A	196	THR
1	A	201	LYS
1	A	217	ASN
1	A	220	GLU
1	A	224	GLU
1	A	249	GLN
2	B	75	LYS
2	B	89	GLN
2	B	97	ARG
3	D	4	GLN

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Mol	Chain	Res	Type
3	D	51	ASN
3	D	59	LYS
3	D	69	SER
3	D	114	GLN
3	D	159	CYS
3	D	161	LEU
3	D	166	MET
3	D	189	ASN
4	E	129	SER
4	E	130	GLU
4	E	144	LEU
4	E	162	LYS
4	E	183	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	HIS
1	A	64	GLN
1	A	90	HIS
1	A	141	GLN
1	A	203	HIS
1	A	217	ASN
2	B	17	ASN
2	B	31	HIS
3	D	5	ASN
3	D	39	GLN
3	D	40	HIS
3	D	51	ASN
3	D	144	ASN
3	D	189	ASN
4	E	11	GLN
4	E	205	HIS
4	E	218	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	$\# Z > 2$
5	JSO	A	301	1	14,14,14	0.97	1 (7%)	16,20,20	0.90	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
5	JSO	A	301	1	-	-	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	301	JSO	C7-C6	2.61	1.45	1.39

There are no bond angle outliers.

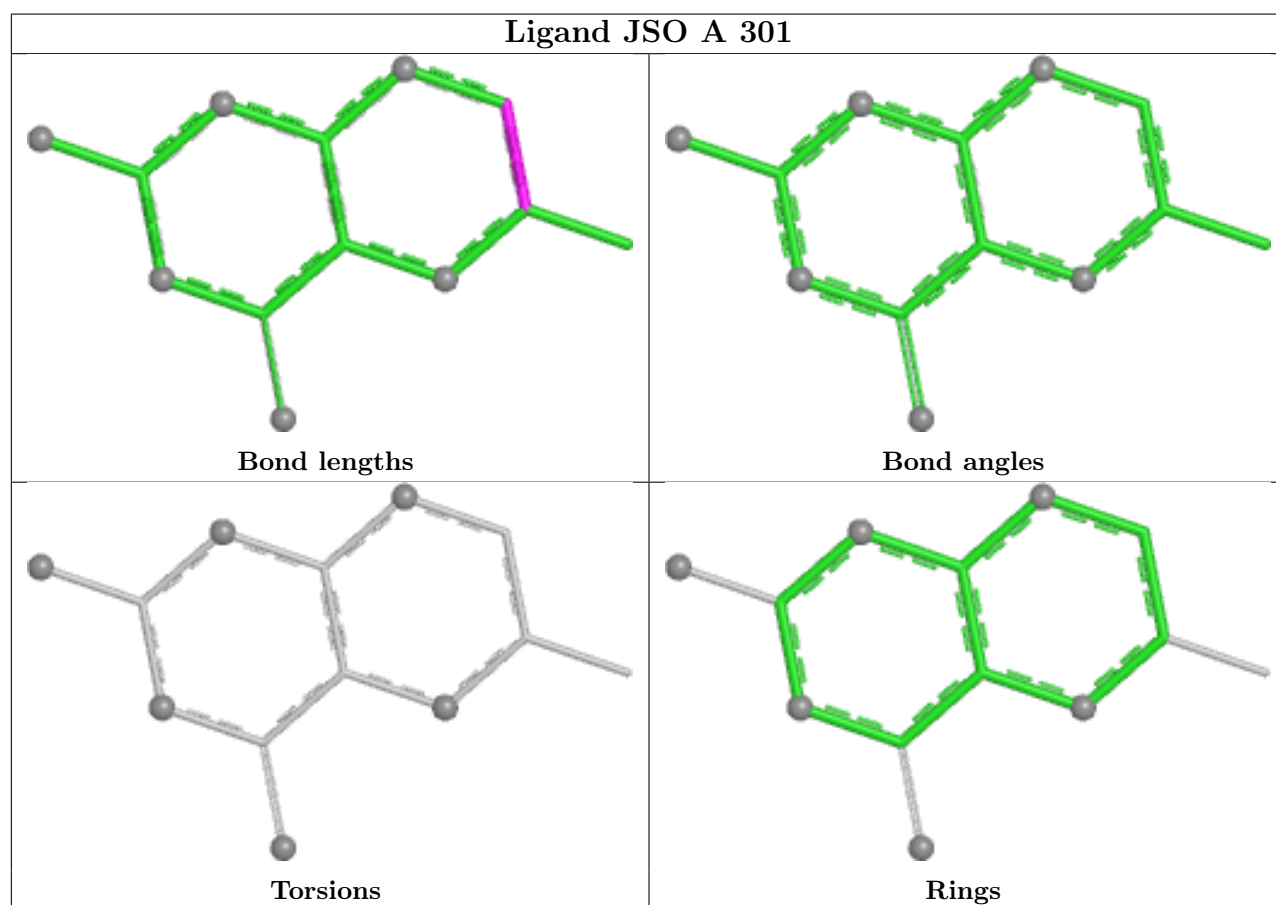
There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	264/290 (91%)	0.88	36 (13%) 7 6	31, 58, 91, 114	1 (0%)
2	B	99/100 (99%)	0.91	7 (7%) 22 21	42, 60, 90, 110	0
3	D	183/205 (89%)	0.75	22 (12%) 9 8	39, 53, 89, 106	0
4	E	242/262 (92%)	0.55	12 (4%) 34 33	26, 49, 73, 86	2 (0%)
All	All	788/857 (91%)	0.75	77 (9%) 13 12	26, 54, 87, 114	3 (0%)

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	264[A]	HIS	5.7
1	A	221	ILE	4.6
1	A	250	SER	4.4
1	A	18	GLY	4.4
1	A	254	TYR	4.3
1	A	265	MET	3.6
3	D	3	GLY	3.6
1	A	132	VAL	3.4
1	A	0	MET	3.4
3	D	127	LYS	3.4
1	A	228	GLY	3.2
3	D	189	ASN	3.2
3	D	180	SER	3.2
1	A	16	ILE	3.1
4	E	162	LYS	3.1
1	A	217	ASN	3.0
4	E	2	ALA	3.0
3	D	124	ARG	3.0
1	A	268	GLN	3.0
1	A	196	THR	2.9
1	A	134	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	19	VAL	2.7
3	D	123	LEU	2.7
3	D	125	ASP	2.7
2	B	14	PRO	2.6
1	A	197	ALA	2.6
4	E	197	THR	2.6
3	D	72	TYR	2.6
3	D	185	ALA	2.6
1	A	246	LEU	2.6
3	D	145	VAL	2.5
1	A	267	LEU	2.5
4	E	1	ASN	2.5
2	B	20	SER	2.5
2	B	97	ARG	2.5
2	B	67	TYR	2.4
1	A	215	MET	2.4
1	A	192	PHE	2.4
3	D	168	PHE	2.4
1	A	244	ILE	2.4
3	D	187	ALA	2.4
3	D	182	PHE	2.4
2	B	49	VAL	2.4
3	D	70	LYS	2.4
3	D	144	ASN	2.3
1	A	189	LYS	2.3
4	E	158	TRP	2.3
2	B	0	MET	2.3
4	E	174	PRO	2.3
4	E	161	GLY	2.3
4	E	3	GLY	2.3
4	E	181	LEU	2.3
3	D	148	SER	2.2
1	A	191	THR	2.2
1	A	222	VAL	2.2
3	D	186	ASN	2.2
1	A	137	HIS	2.2
1	A	266	VAL	2.1
1	A	185	ARG	2.1
1	A	187	ASN	2.1
3	D	166	MET	2.1
3	D	188	PHE	2.1
1	A	17	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	203	HIS	2.1
1	A	263	VAL	2.1
3	D	4	GLN	2.1
1	A	195	VAL	2.1
3	D	69	SER	2.1
1	A	249	GLN	2.1
4	E	165	HIS	2.1
1	A	199	PHE	2.1
3	D	29	SER	2.1
4	E	241	ALA	2.1
4	E	116	ASN	2.0
2	B	98	ASP	2.0
1	A	130	LEU	2.0
1	A	198	LEU	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 [i](#)

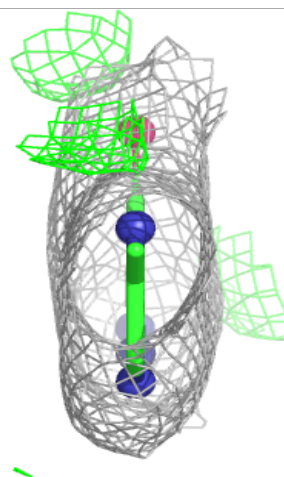
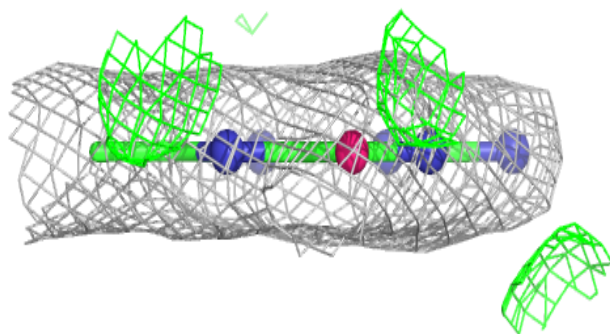
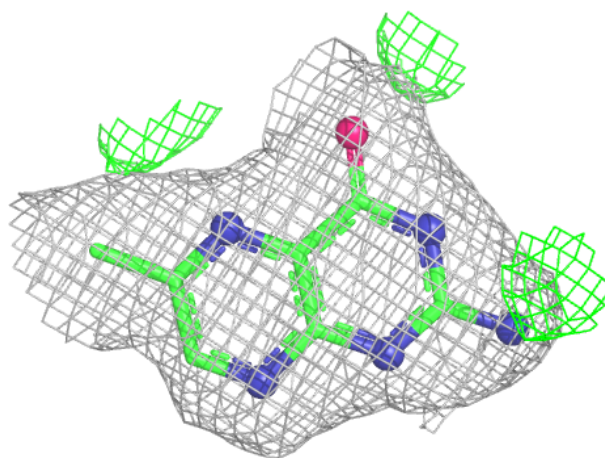
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	JSO	A	301	13/13	0.94	0.08	40,47,51,51	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 JSO 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.