



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:53 AM UTC

PDB ID : 7ZXS / pdb_00007zxs
Title : Crystal structure of DPP9 in complex with a 4-oxo-b-lactam based inhibitor, A295
Authors : Ross, B.; Huber, R.
Deposited on : 2022-05-22
Resolution : 1.81 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

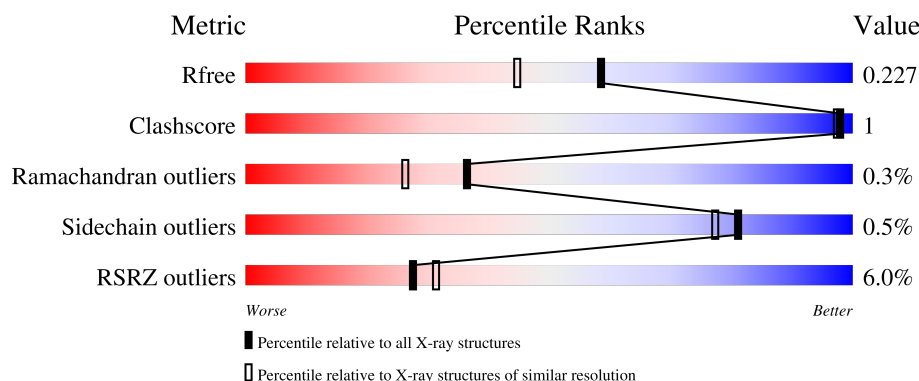
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

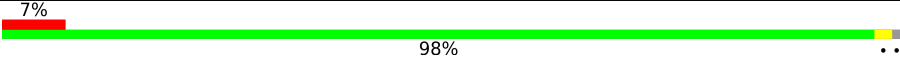
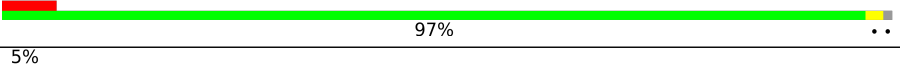
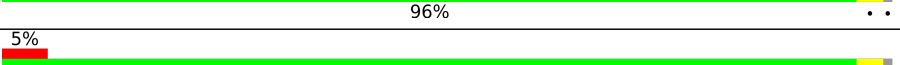
The reported resolution of this entry is 1.81 Å.

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	1112 (1.82-1.82)
Clashscore	190562	1148 (1.82-1.82)
Ramachandran outliers	187476	1140 (1.82-1.82)
Sidechain outliers	187428	1140 (1.82-1.82)
RSRZ outliers	180081	1112 (1.82-1.82)

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	850	 7% 98% ..
1	C	850	 6% 97% ..
2	B	850	 5% 96% ..
2	D	850	 5% 96% ..

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 29394 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 Dipeptidyl peptidase 9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	844	Total	C	N	O	S	145	4	0
			6854	4399	1170	1256	29			
1	C	843	Total	C	N	O	S	155	8	0
			6881	4415	1178	1258	30			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	864	HIS	-	expression tag	UNP Q86TI2
A	865	HIS	-	expression tag	UNP Q86TI2
A	866	HIS	-	expression tag	UNP Q86TI2
A	867	HIS	-	expression tag	UNP Q86TI2
A	868	HIS	-	expression tag	UNP Q86TI2
A	869	HIS	-	expression tag	UNP Q86TI2
C	864	HIS	-	expression tag	UNP Q86TI2
C	865	HIS	-	expression tag	UNP Q86TI2
C	866	HIS	-	expression tag	UNP Q86TI2
C	867	HIS	-	expression tag	UNP Q86TI2
C	868	HIS	-	expression tag	UNP Q86TI2
C	869	HIS	-	expression tag	UNP Q86TI2

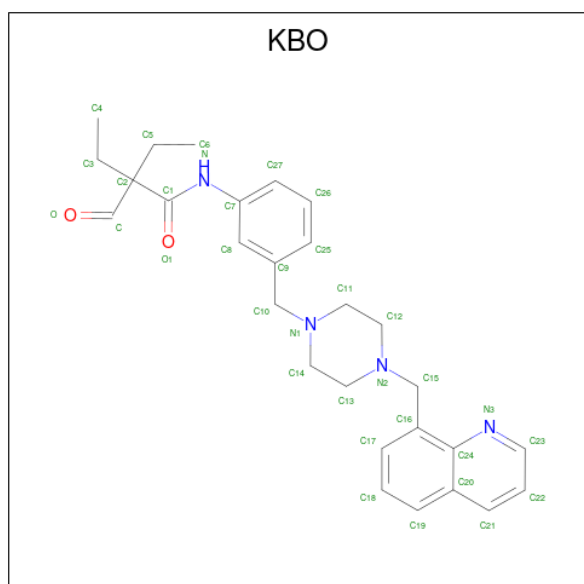
- Molecule 2 is a protein called Dipeptidyl peptidase 9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	839	Total	C	N	O	S	127	10	0
			6855	4396	1174	1254	31			
2	D	838	Total	C	N	O	S	122	8	0
			6830	4379	1168	1252	31			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	864	HIS	-	expression tag	UNP Q86TI2
B	865	HIS	-	expression tag	UNP Q86TI2
B	866	HIS	-	expression tag	UNP Q86TI2
B	867	HIS	-	expression tag	UNP Q86TI2
B	868	HIS	-	expression tag	UNP Q86TI2
B	869	HIS	-	expression tag	UNP Q86TI2
D	864	HIS	-	expression tag	UNP Q86TI2
D	865	HIS	-	expression tag	UNP Q86TI2
D	866	HIS	-	expression tag	UNP Q86TI2
D	867	HIS	-	expression tag	UNP Q86TI2
D	868	HIS	-	expression tag	UNP Q86TI2
D	869	HIS	-	expression tag	UNP Q86TI2

- Molecule 3 is 2-ethyl-2-methanoyl- {N}-[3-[[4-(quinolin-8-ylmethyl)piperazin-1-yl]methyl]phenyl]butanamide (CCD ID: KBO) (formula: $C_{28}H_{34}N_4O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	10	0
			34	28	4	2		
3	B	1	Total	C	N	O	10	0
			34	28	4	2		
3	C	1	Total	C	N	O	10	0
			34	28	4	2		
3	D	1	Total	C	N	O	10	0
			34	28	4	2		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

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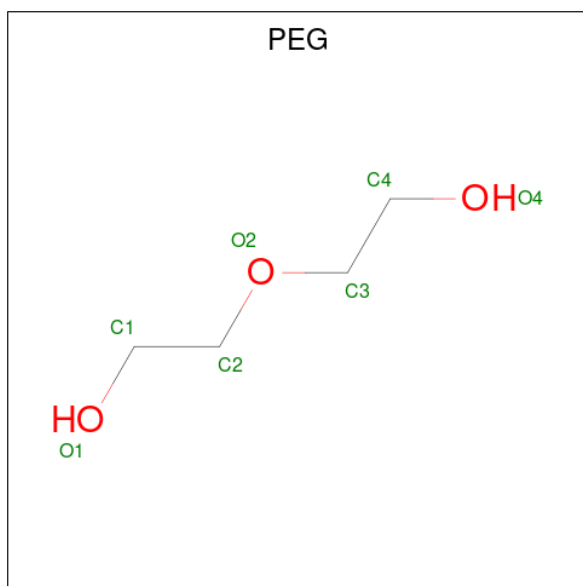
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0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

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

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			7	4	3		
5	C	1	Total	C	O	0	0
			7	4	3		
5	D	1	Total	C	O	0	0
			7	4	3		

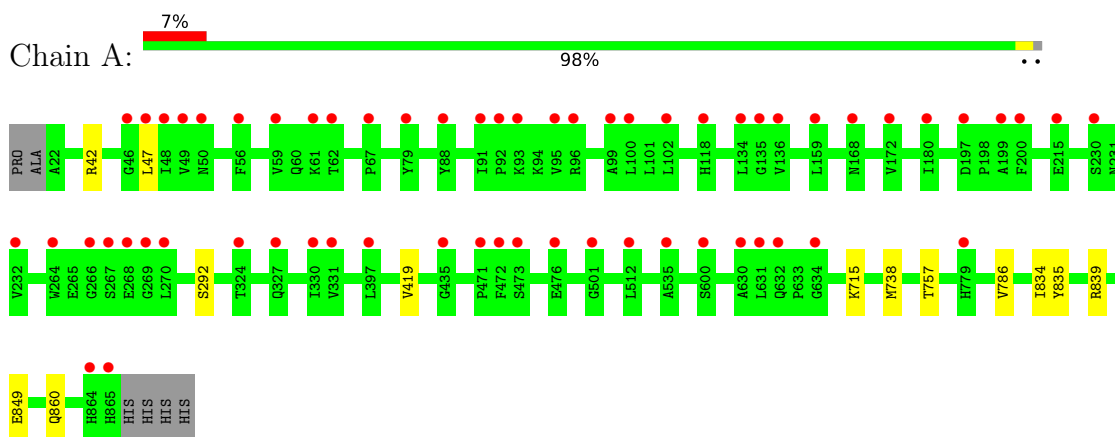
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	331	Total 331	O 331	0	0
6	B	425	Total 425	O 425	0	0
6	C	419	Total 419	O 419	0	0
6	D	478	Total 478	O 478	0	0

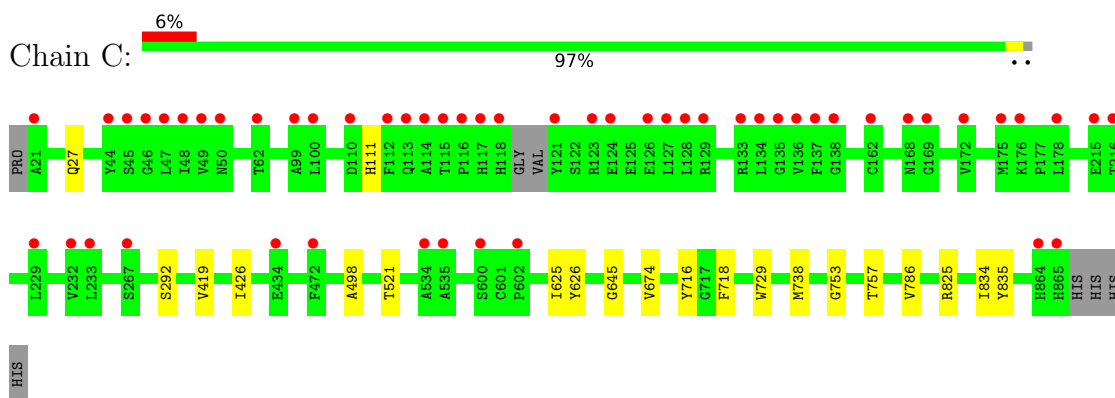
3 Residue-property plots [i](#)

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.

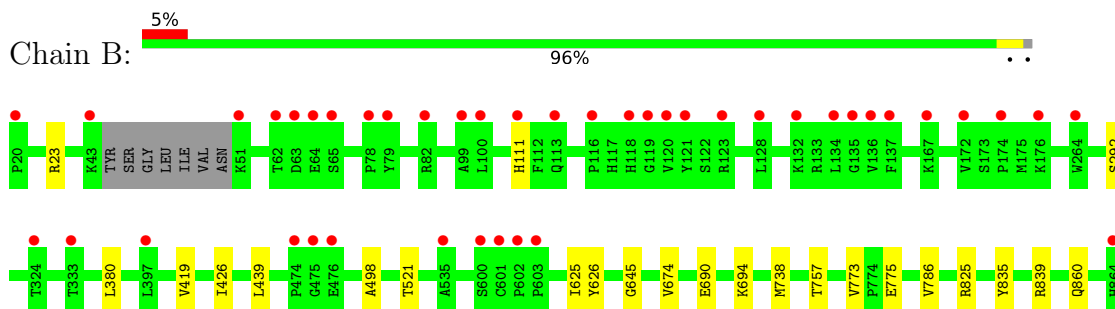
• Molecule 1: Dipeptidyl peptidase 9



• Molecule 1: Dipeptidyl peptidase 9

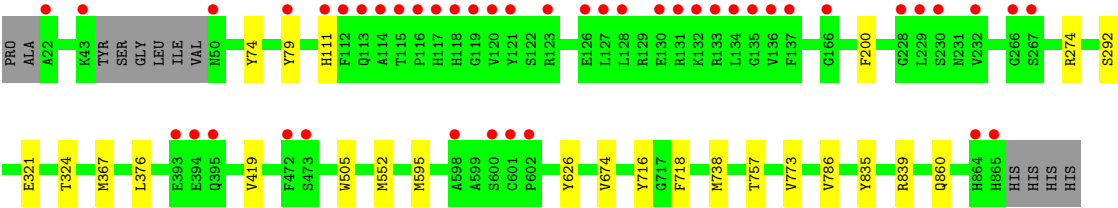


• Molecule 2: Dipeptidyl peptidase 9



H965
HIS
HIS
HIS
HIS

• Molecule 2: Dipeptidyl peptidase 9



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	88.35Å 106.18Å 121.13Å 65.25° 70.43° 75.86°	Depositor
Resolution (Å)	106.09 – 1.81 106.09 – 1.81	Depositor EDS
% Data completeness (in resolution range)	64.6 (106.09-1.81) 64.8 (106.09-1.81)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 1.81Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.175 , 0.203 (Not available) , 0.227	Depositor DCC
R_{free} test set	1707 reflections (0.74%)	wwPDB-VP
Wilson B-factor (Å ²)	30.1	Xtriage
Anisotropy	0.023	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 32.6	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	0.96	EDS
Total number of atoms	29394	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.27% 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: PEG, EDO, KBO, CSO

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	0.99	0/7053	1.26	1/9571 (0.0%)
1	C	0.97	0/7080	1.25	1/9605 (0.0%)
2	B	0.97	0/7061	1.26	1/9581 (0.0%)
2	D	0.97	0/7035	1.25	2/9547 (0.0%)
All	All	0.97	0/28229	1.26	5/38304 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	835	TYR	CB-CA-C	7.55	118.50	110.65
1	A	835	TYR	CB-CA-C	6.90	117.83	110.65
2	D	835	TYR	CB-CA-C	6.66	117.58	110.65
1	C	835	TYR	CB-CA-C	6.37	117.27	110.65
2	D	79	TYR	CB-CA-C	5.26	118.31	109.89

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	6854	0	6625	5	0
1	C	6881	0	6650	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	6855	0	6623	14	0
2	D	6830	0	6594	11	0
3	A	34	0	0	0	0
3	B	34	0	0	0	0
3	C	34	0	0	0	0
3	D	34	0	0	0	0
4	A	28	0	42	0	0
4	B	48	0	72	0	0
4	C	40	0	60	0	0
4	D	48	0	72	0	0
5	B	7	0	10	0	0
5	C	7	0	10	0	0
5	D	7	0	10	0	0
6	A	331	0	0	1	0
6	B	425	0	0	3	0
6	C	419	0	0	1	0
6	D	478	0	0	1	0
All	All	29394	0	26768	37	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 1.

All (37) 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:825[B]:ARG:HH11	2:B:825[B]:ARG:HG2	1.44	0.81
2:B:825[A]:ARG:HH11	2:B:825[A]:ARG:HG3	1.59	0.67
2:B:825[B]:ARG:HG2	2:B:825[B]:ARG:NH1	2.06	0.64
2:B:738:MET:HE3	2:B:786:VAL:HG12	1.81	0.62
2:B:825[A]:ARG:HG3	2:B:825[A]:ARG:NH1	2.19	0.55
2:D:738:MET:HE3	2:D:786:VAL:HG12	1.89	0.55
1:A:834[B]:ILE:HD11	1:C:834[B]:ILE:HD11	1.90	0.54
1:C:645:GLY:HA2	6:C:1171:HOH:O	2.09	0.53
2:D:200:PHE:CZ	2:D:324:THR:HG21	2.44	0.52
1:C:716:TYR:HB3	1:C:718:PHE:CE2	2.46	0.50
2:D:716:TYR:HB3	2:D:718:PHE:CE2	2.47	0.50
1:C:757:THR:HA	1:C:786:VAL:HG22	1.95	0.49
2:B:690:GLU:HG3	6:B:1003:HOH:O	2.12	0.49
2:D:367:MET:HE3	2:D:376:LEU:HD11	1.95	0.47
2:D:757:THR:HA	2:D:786:VAL:HG22	1.96	0.47
2:B:380:LEU:HD11	2:B:439:LEU:HD22	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:738:MET:HE3	1:A:786:VAL:HG12	1.98	0.46
2:B:426:ILE:HD12	2:B:498:ALA:HB2	1.98	0.45
2:B:773:VAL:HG23	2:B:775:GLU:HG2	1.99	0.45
2:B:23:ARG:HA	2:B:23:ARG:HD3	1.84	0.45
1:C:626:TYR:HB2	1:C:674:VAL:HB	1.98	0.44
2:B:645:GLY:HA2	6:B:1118:HOH:O	2.17	0.44
2:B:694:LYS:NZ	6:B:1014:HOH:O	2.50	0.44
2:D:74:TYR:CZ	2:D:595:MET:HE1	2.52	0.44
1:A:757:THR:HA	1:A:786:VAL:HG22	2.00	0.43
2:B:757:THR:HA	2:B:786:VAL:HG22	2.00	0.43
2:D:505:TRP:CD1	2:D:552:MET:HB3	2.53	0.43
1:A:715:LYS:NZ	6:A:1005:HOH:O	2.51	0.43
2:D:274:ARG:HG2	2:D:321:GLU:HG2	2.01	0.43
1:C:738:MET:HE3	1:C:786:VAL:HG12	2.01	0.42
1:C:729:TRP:CE3	1:C:753:GLY:HA3	2.54	0.42
2:B:626:TYR:HB2	2:B:674:VAL:HB	2.01	0.42
1:A:42:ARG:HD3	1:A:849:GLU:HG2	2.02	0.41
2:D:626:TYR:HB2	2:D:674:VAL:HB	2.02	0.41
2:D:773:VAL:HG13	6:D:1380:HOH:O	2.19	0.41
2:D:738:MET:CE	2:D:786:VAL:HG12	2.51	0.40
1:C:426:ILE:HD12	1:C:498:ALA:HB2	2.03	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	845/850 (99%)	821 (97%)	22 (3%)	2 (0%)	43	33
1	C	846/850 (100%)	823 (97%)	21 (2%)	2 (0%)	43	33
2	B	845/850 (99%)	821 (97%)	21 (2%)	3 (0%)	30	19
2	D	842/850 (99%)	824 (98%)	15 (2%)	3 (0%)	30	19

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	3378/3400 (99%)	3289 (97%)	79 (2%)	10 (0%)	36	26

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	839	ARG
1	C	111	HIS
2	D	111	HIS
2	B	111	HIS
2	B	839	ARG
2	D	839	ARG
1	A	419	VAL
2	B	419	VAL
1	C	419	VAL
2	D	419	VAL

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	745/746 (100%)	742 (100%)	3 (0%)	84	81
1	C	748/746 (100%)	743 (99%)	5 (1%)	76	70
2	B	747/747 (100%)	743 (100%)	4 (0%)	81	77
2	D	745/747 (100%)	742 (100%)	3 (0%)	84	81
All	All	2985/2986 (100%)	2970 (100%)	15 (0%)	81	77

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	LEU
1	A	292	SER
1	A	860	GLN
2	B	292	SER
2	B	521	THR

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Mol	Chain	Res	Type
2	B	625	ILE
2	B	860	GLN
1	C	27	GLN
1	C	292	SER
1	C	521	THR
1	C	625	ILE
1	C	825	ARG
2	D	292[A]	SER
2	D	292[B]	SER
2	D	860	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	GLN
1	A	161	HIS
1	A	614	HIS
2	B	161	HIS
2	B	226	HIS
2	B	247	GLN
2	B	323	GLN
2	B	798	ASN
2	B	831	GLN
1	C	29	HIS
1	C	161	HIS
1	C	226	HIS
1	C	432	GLN
1	C	614	HIS
1	C	810	ASN
2	D	161	HIS
2	D	555	ASN
2	D	865	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	CSO	C	844	1	3,6,7	0.73	0	1,6,8	0.09	0
1	CSO	A	844	1	3,6,7	0.60	0	1,6,8	0.26	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
1	CSO	C	844	1	-	0/1/5/7	-
1	CSO	A	844	1	-	0/1/5/7	-

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.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

48 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	$\# Z > 2$
4	EDO	D	911	-	3,3,3	0.24	0	2,2,2	0.34	0
4	EDO	B	906	-	3,3,3	0.13	0	2,2,2	0.28	0
4	EDO	A	902	-	3,3,3	0.12	0	2,2,2	0.26	0
4	EDO	B	902	-	3,3,3	0.11	0	2,2,2	0.29	0
4	EDO	B	914	-	3,3,3	0.11	0	2,2,2	0.37	0
4	EDO	D	914	-	3,3,3	0.08	0	2,2,2	0.19	0
4	EDO	A	905	-	3,3,3	0.09	0	2,2,2	0.20	0
4	EDO	A	908	-	3,3,3	0.20	0	2,2,2	0.07	0
4	EDO	B	904	-	3,3,3	0.07	0	2,2,2	0.30	0
4	EDO	B	905	-	3,3,3	0.07	0	2,2,2	0.18	0
4	EDO	B	910	-	3,3,3	0.05	0	2,2,2	0.13	0
4	EDO	D	902	-	3,3,3	0.01	0	2,2,2	0.05	0
4	EDO	A	907	-	3,3,3	0.20	0	2,2,2	0.18	0
3	KBO	A	901	1	35,37,37	0.93	0	45,51,51	1.48	9 (20%)
4	EDO	C	904	-	3,3,3	0.18	0	2,2,2	0.27	0
4	EDO	C	905	-	3,3,3	0.04	0	2,2,2	0.19	0
5	PEG	B	913	-	6,6,6	0.17	0	5,5,5	0.13	0
3	KBO	D	901	2	35,37,37	0.96	1 (2%)	45,51,51	1.51	8 (17%)
5	PEG	D	912	-	6,6,6	0.18	0	5,5,5	0.15	0
4	EDO	D	908	-	3,3,3	0.13	0	2,2,2	0.16	0
4	EDO	D	903	-	3,3,3	0.14	0	2,2,2	0.24	0
4	EDO	B	903	-	3,3,3	0.07	0	2,2,2	0.07	0
5	PEG	C	909	-	6,6,6	0.19	0	5,5,5	0.10	0
4	EDO	B	907	-	3,3,3	0.19	0	2,2,2	0.34	0
4	EDO	D	909	-	3,3,3	0.11	0	2,2,2	0.37	0
4	EDO	C	908	-	3,3,3	0.18	0	2,2,2	0.33	0
4	EDO	C	906	-	3,3,3	0.09	0	2,2,2	0.05	0
4	EDO	C	907	-	3,3,3	0.06	0	2,2,2	0.21	0
4	EDO	C	912	-	3,3,3	0.07	0	2,2,2	0.03	0
4	EDO	D	913	-	3,3,3	0.16	0	2,2,2	0.45	0
4	EDO	B	912	-	3,3,3	0.23	0	2,2,2	0.26	0
4	EDO	B	908	-	3,3,3	0.11	0	2,2,2	0.11	0
4	EDO	C	902	-	3,3,3	0.04	0	2,2,2	0.02	0
4	EDO	D	906	-	3,3,3	0.11	0	2,2,2	0.28	0
4	EDO	C	911	-	3,3,3	0.07	0	2,2,2	0.24	0
4	EDO	C	903	-	3,3,3	0.10	0	2,2,2	0.20	0
4	EDO	C	910	-	3,3,3	0.10	0	2,2,2	0.30	0
3	KBO	B	901	2	35,37,37	0.99	1 (2%)	45,51,51	1.67	9 (20%)
4	EDO	A	904	-	3,3,3	0.07	0	2,2,2	0.23	0
4	EDO	D	904	-	3,3,3	0.15	0	2,2,2	0.26	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	D	905	-	3,3,3	0.34	0	2,2,2	0.47	0
4	EDO	D	910	-	3,3,3	0.17	0	2,2,2	0.11	0
4	EDO	B	911	-	3,3,3	0.05	0	2,2,2	0.18	0
4	EDO	A	903	-	3,3,3	0.12	0	2,2,2	0.06	0
4	EDO	A	906	-	3,3,3	0.10	0	2,2,2	0.21	0
4	EDO	B	909	-	3,3,3	0.18	0	2,2,2	0.40	0
3	KBO	C	901	1	35,37,37	0.96	1 (2%)	45,51,51	1.59	8 (17%)
4	EDO	D	907	-	3,3,3	0.09	0	2,2,2	0.31	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
4	EDO	D	911	-	-	1/1/1/1	-
4	EDO	B	906	-	-	0/1/1/1	-
4	EDO	A	902	-	-	0/1/1/1	-
4	EDO	B	902	-	-	1/1/1/1	-
4	EDO	B	914	-	-	0/1/1/1	-
4	EDO	D	914	-	-	1/1/1/1	-
4	EDO	A	905	-	-	0/1/1/1	-
4	EDO	A	908	-	-	0/1/1/1	-
4	EDO	B	904	-	-	1/1/1/1	-
4	EDO	B	905	-	-	0/1/1/1	-
4	EDO	B	910	-	-	1/1/1/1	-
4	EDO	D	902	-	-	0/1/1/1	-
4	EDO	A	907	-	-	1/1/1/1	-
3	KBO	A	901	1	-	5/24/37/37	0/4/4/4
4	EDO	C	904	-	-	0/1/1/1	-
4	EDO	C	905	-	-	0/1/1/1	-
5	PEG	B	913	-	-	1/4/4/4	-
3	KBO	D	901	2	-	5/24/37/37	0/4/4/4
5	PEG	D	912	-	-	2/4/4/4	-
4	EDO	D	908	-	-	1/1/1/1	-
4	EDO	D	903	-	-	0/1/1/1	-
4	EDO	B	903	-	-	0/1/1/1	-
5	PEG	C	909	-	-	0/4/4/4	-
4	EDO	B	907	-	-	1/1/1/1	-
4	EDO	D	909	-	-	0/1/1/1	-
4	EDO	C	908	-	-	1/1/1/1	-
4	EDO	C	906	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	C	907	-	-	1/1/1/1	-
4	EDO	C	912	-	-	0/1/1/1	-
4	EDO	D	913	-	-	1/1/1/1	-
4	EDO	B	912	-	-	0/1/1/1	-
4	EDO	B	908	-	-	0/1/1/1	-
4	EDO	C	902	-	-	1/1/1/1	-
4	EDO	D	906	-	-	1/1/1/1	-
4	EDO	C	911	-	-	0/1/1/1	-
4	EDO	C	903	-	-	0/1/1/1	-
4	EDO	C	910	-	-	0/1/1/1	-
3	KBO	B	901	2	-	3/24/37/37	0/4/4/4
4	EDO	A	904	-	-	1/1/1/1	-
4	EDO	D	904	-	-	0/1/1/1	-
4	EDO	D	905	-	-	0/1/1/1	-
4	EDO	D	910	-	-	0/1/1/1	-
4	EDO	B	911	-	-	0/1/1/1	-
4	EDO	A	903	-	-	0/1/1/1	-
4	EDO	A	906	-	-	1/1/1/1	-
4	EDO	B	909	-	-	1/1/1/1	-
3	KBO	C	901	1	-	8/24/37/37	0/4/4/4
4	EDO	D	907	-	-	1/1/1/1	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	901	KBO	C8-C7	2.21	1.43	1.39
3	D	901	KBO	C8-C7	2.19	1.43	1.39
3	C	901	KBO	C8-C7	2.04	1.42	1.39

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	901	KBO	C16-C15-N2	-5.64	106.11	114.07
3	C	901	KBO	C16-C15-N2	-4.63	107.54	114.07
3	D	901	KBO	C16-C15-N2	-4.26	108.05	114.07
3	C	901	KBO	C23-N3-C24	4.05	122.08	117.31
3	D	901	KBO	C23-N3-C24	4.05	122.08	117.31
3	B	901	KBO	C23-N3-C24	4.01	122.03	117.31
3	A	901	KBO	C23-N3-C24	3.96	121.98	117.31
3	C	901	KBO	C2-C1-N	3.90	120.83	115.68
3	B	901	KBO	C2-C1-N	3.86	120.77	115.68
3	D	901	KBO	C2-C1-N	3.44	120.22	115.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	KBO	C2-C1-N	3.38	120.14	115.68
3	D	901	KBO	C20-C24-N3	-3.37	119.07	122.63
3	C	901	KBO	C20-C24-N3	-3.36	119.08	122.63
3	A	901	KBO	C20-C24-N3	-3.28	119.17	122.63
3	B	901	KBO	C20-C24-N3	-3.13	119.32	122.63
3	A	901	KBO	C16-C15-N2	-2.96	109.88	114.07
3	C	901	KBO	C27-C7-C8	-2.95	116.09	119.66
3	B	901	KBO	C26-C27-C7	2.74	122.88	119.73
3	B	901	KBO	C27-C7-C8	-2.68	116.42	119.66
3	D	901	KBO	C27-C7-C8	-2.59	116.52	119.66
3	A	901	KBO	C27-C7-C8	-2.55	116.57	119.66
3	B	901	KBO	C22-C23-N3	-2.52	120.28	123.97
3	C	901	KBO	C22-C23-N3	-2.48	120.32	123.97
3	D	901	KBO	C22-C23-N3	-2.44	120.39	123.97
3	A	901	KBO	C22-C23-N3	-2.43	120.40	123.97
3	A	901	KBO	C15-C16-C24	2.26	122.11	118.59
3	C	901	KBO	C26-C27-C7	2.26	122.33	119.73
3	D	901	KBO	C16-C24-N3	2.22	121.49	118.16
3	C	901	KBO	C16-C24-N3	2.19	121.45	118.16
3	D	901	KBO	C15-C16-C24	2.19	121.99	118.59
3	A	901	KBO	C16-C24-N3	2.17	121.42	118.16
3	A	901	KBO	C8-C7-N	2.09	126.94	120.13
3	B	901	KBO	C16-C24-N3	2.09	121.30	118.16
3	B	901	KBO	C25-C9-C8	2.09	121.42	118.55

There are no chirality outliers.

All (41) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	901	KBO	O-C-C2-C3
3	B	901	KBO	O-C-C2-C3
3	D	901	KBO	O-C-C2-C3
3	D	901	KBO	C16-C15-N2-C12
3	D	901	KBO	C16-C15-N2-C13
4	D	906	EDO	O1-C1-C2-O2
5	D	912	PEG	O1-C1-C2-O2
3	C	901	KBO	C16-C15-N2-C12
3	C	901	KBO	C16-C15-N2-C13
3	C	901	KBO	C3-C2-C5-C6
4	A	907	EDO	O1-C1-C2-O2
4	D	907	EDO	O1-C1-C2-O2
4	D	911	EDO	O1-C1-C2-O2

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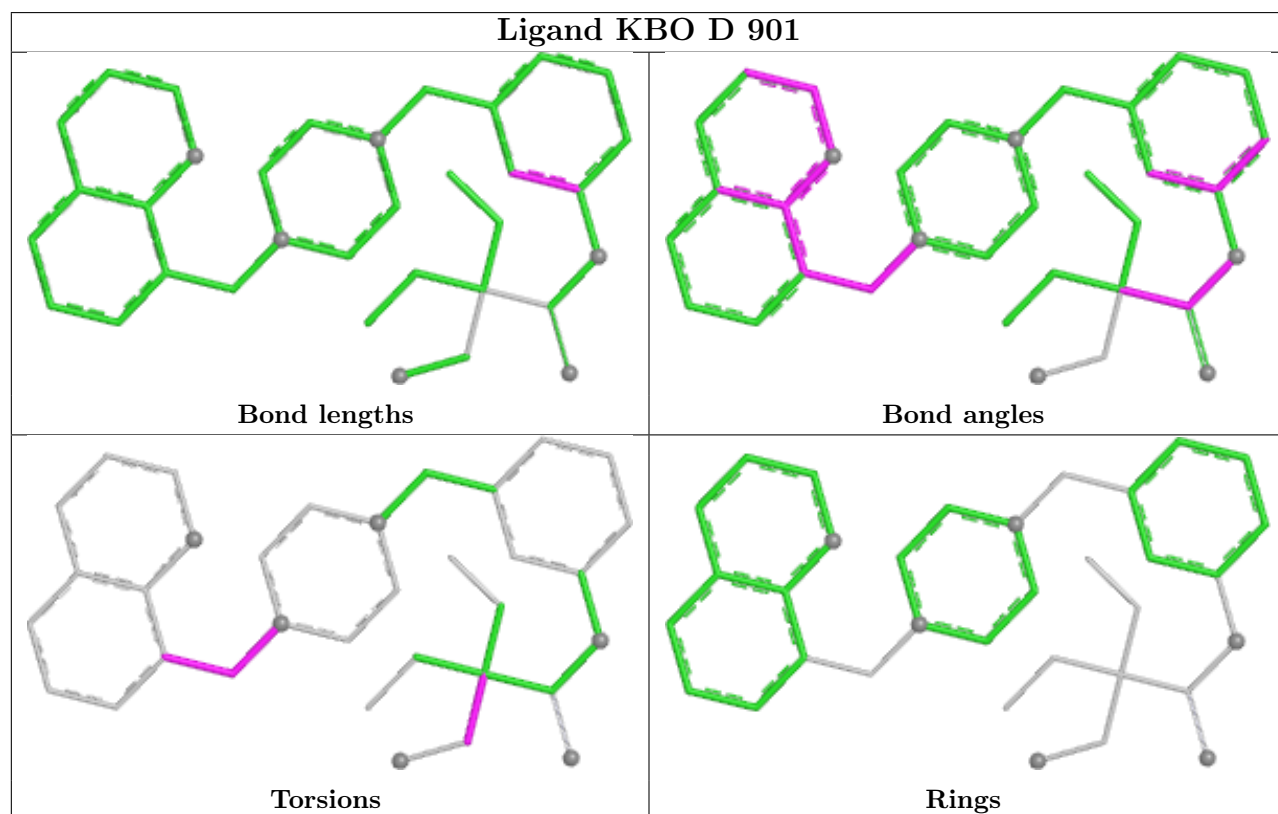
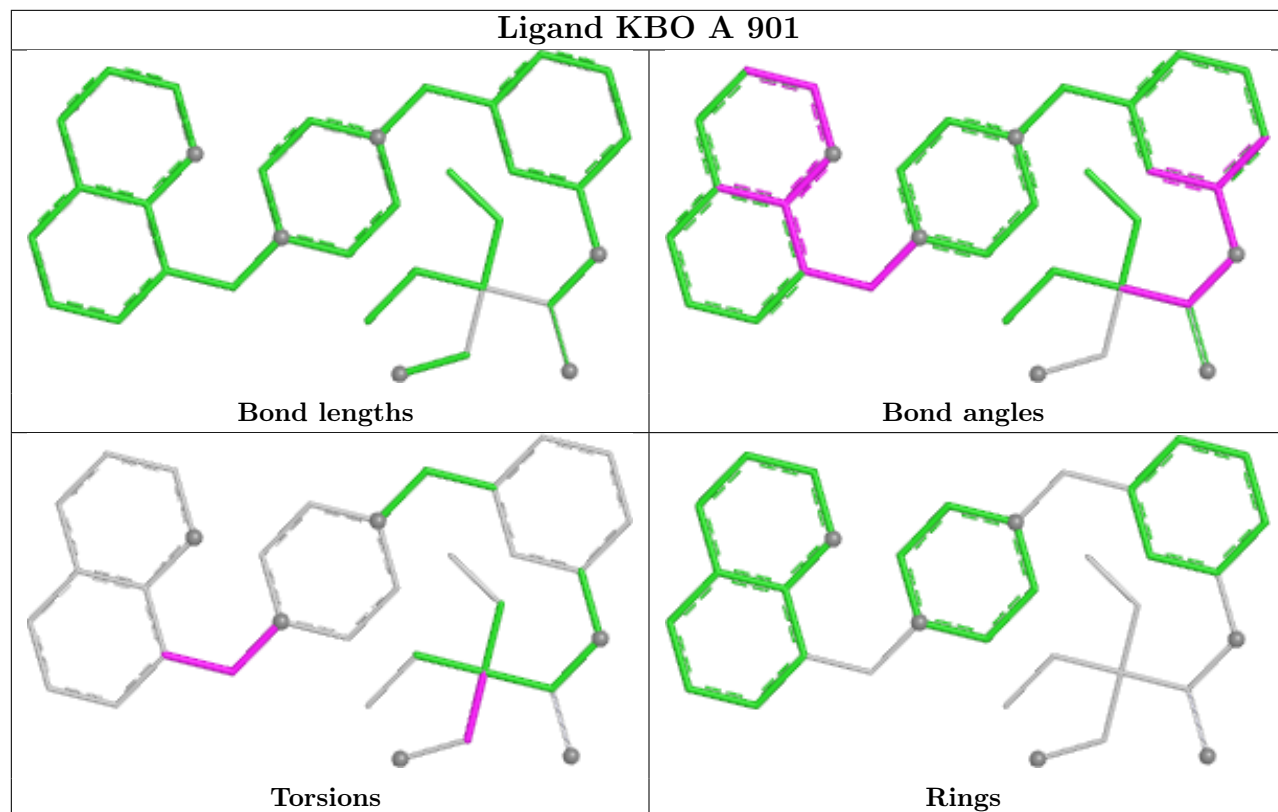
Mol	Chain	Res	Type	Atoms
4	D	914	EDO	O1-C1-C2-O2
3	A	901	KBO	C16-C15-N2-C12
4	B	902	EDO	O1-C1-C2-O2
4	C	902	EDO	O1-C1-C2-O2
4	D	913	EDO	O1-C1-C2-O2
5	B	913	PEG	O2-C3-C4-O4
3	A	901	KBO	N2-C15-C16-C17
3	A	901	KBO	C16-C15-N2-C13
3	D	901	KBO	N2-C15-C16-C17
3	B	901	KBO	N2-C15-C16-C17
3	C	901	KBO	N2-C15-C16-C17
5	D	912	PEG	C4-C3-O2-C2
3	C	901	KBO	O-C-C2-C3
3	C	901	KBO	O-C-C2-C5
3	A	901	KBO	N2-C15-C16-C24
3	C	901	KBO	N2-C15-C16-C24
3	D	901	KBO	N2-C15-C16-C24
4	B	910	EDO	O1-C1-C2-O2
4	B	907	EDO	O1-C1-C2-O2
4	B	909	EDO	O1-C1-C2-O2
4	C	907	EDO	O1-C1-C2-O2
3	B	901	KBO	N2-C15-C16-C24
4	A	904	EDO	O1-C1-C2-O2
4	A	906	EDO	O1-C1-C2-O2
4	B	904	EDO	O1-C1-C2-O2
4	C	908	EDO	O1-C1-C2-O2
4	D	908	EDO	O1-C1-C2-O2
3	C	901	KBO	C1-C2-C5-C6

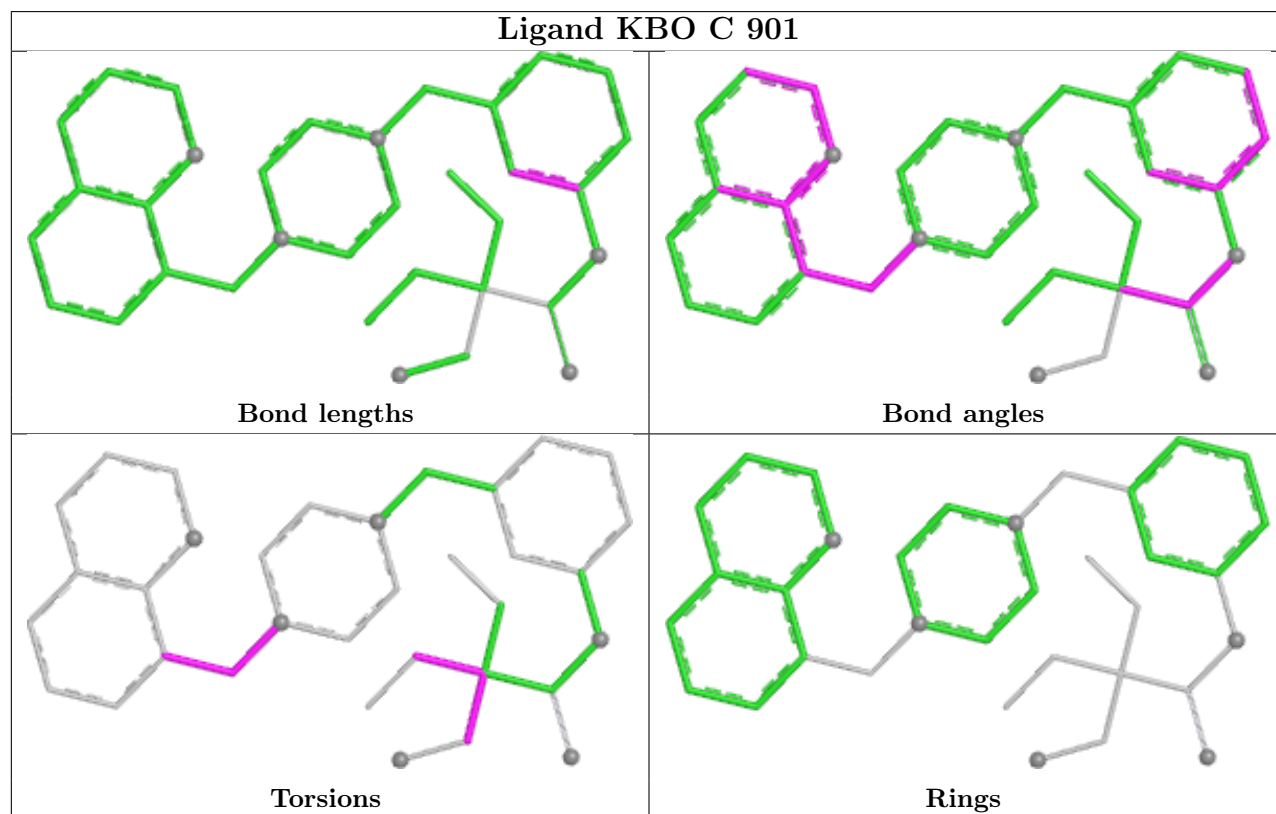
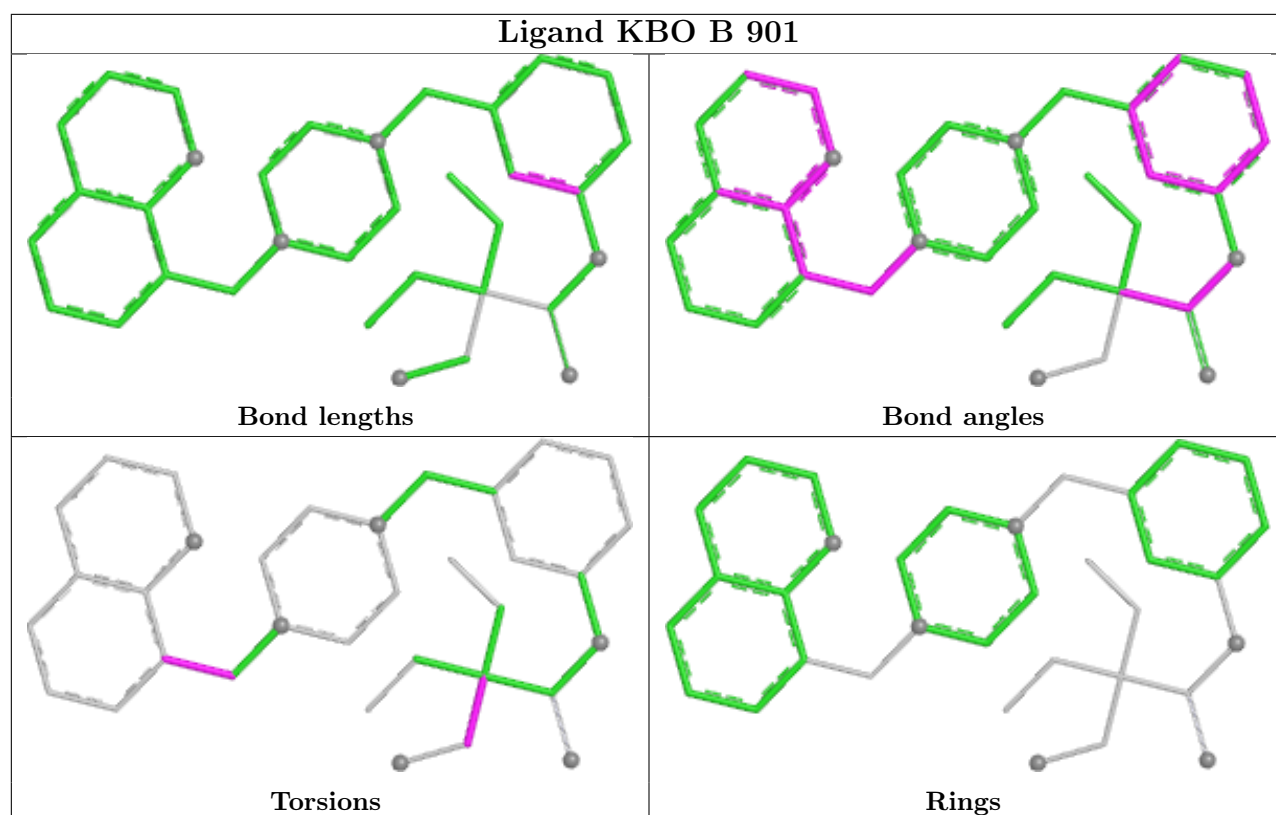
There are no ring outliers.

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	843/850 (99%)	0.52	61 (7%)	21	24	11, 40, 74, 107	61 (7%)
1	C	842/850 (99%)	0.19	53 (6%)	26	30	11, 31, 64, 104	63 (7%)
2	B	839/850 (98%)	0.13	43 (5%)	33	38	10, 30, 63, 92	56 (6%)
2	D	838/850 (98%)	0.12	45 (5%)	31	35	11, 30, 65, 110	51 (6%)
All	All	3362/3400 (98%)	0.24	202 (6%)	27	31	10, 33, 68, 110	231 (6%)

All (202) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	136	VAL	5.1
1	A	47	LEU	4.9
2	B	79	TYR	4.8
1	C	47	LEU	4.7
2	B	602	PRO	4.6
2	D	602	PRO	4.6
1	A	48	ILE	4.5
1	C	21	ALA	4.4
2	D	79	TYR	4.3
1	C	135	GLY	4.3
2	D	121	TYR	4.2
1	A	49	VAL	4.2
2	D	116	PRO	4.2
1	A	62	THR	4.2
2	B	172	VAL	4.1
2	D	120	VAL	4.1
1	A	270	LEU	4.1
1	C	134	LEU	4.0
1	C	136	VAL	4.0
2	D	117	HIS	4.0
1	A	50	ASN	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	99	ALA	3.9
1	C	114	ALA	3.9
1	A	61	LYS	3.8
2	D	135	GLY	3.7
1	C	172	VAL	3.7
2	B	51	LYS	3.7
2	B	65	SER	3.7
2	D	134	LEU	3.6
1	A	435	GLY	3.6
1	C	48	ILE	3.6
2	B	118	HIS	3.6
1	A	99	ALA	3.5
1	C	600	SER	3.5
1	A	269	GLY	3.5
2	B	113	GLN	3.4
1	A	79	TYR	3.4
1	C	121	TYR	3.4
2	B	120	VAL	3.4
2	D	22	ALA	3.4
2	D	864	HIS	3.4
1	C	49	VAL	3.3
1	C	169	GLY	3.3
1	A	266	GLY	3.3
1	C	45	SER	3.3
1	A	331	VAL	3.2
1	C	128	LEU	3.2
2	D	232	VAL	3.1
2	D	600	SER	3.1
1	C	113	GLN	3.1
1	C	116	PRO	3.1
2	D	128	LEU	3.1
1	A	535	ALA	3.1
2	D	118	HIS	3.1
1	A	95	VAL	3.1
1	C	44	TYR	3.0
1	A	264	TRP	3.0
2	B	600	SER	3.0
2	D	112	PHE	3.0
2	D	123	ARG	3.0
2	D	229	LEU	3.0
1	A	267	SER	3.0
1	C	112	PHE	3.0

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Mol	Chain	Res	Type	RSRZ
2	B	603	PRO	3.0
2	B	535	ALA	3.0
2	B	121	TYR	2.9
2	D	113	GLN	2.9
2	D	131	ARG	2.9
1	A	472	PHE	2.9
1	C	46	GLY	2.9
1	C	124	GLU	2.9
1	C	602	PRO	2.9
1	C	168	ASN	2.9
2	D	137	PHE	2.9
2	D	114	ALA	2.9
2	D	111	HIS	2.9
2	B	136	VAL	2.9
1	A	67	PRO	2.9
1	C	535	ALA	2.9
1	C	162	CYS	2.9
2	D	127	LEU	2.9
1	A	46	GLY	2.9
2	D	266	GLY	2.9
2	B	99	ALA	2.8
2	B	476	GLU	2.8
1	C	175	MET	2.8
1	A	134	LEU	2.8
1	A	330	ILE	2.8
2	D	119	GLY	2.8
1	C	137	PHE	2.7
2	D	472	PHE	2.7
1	A	100	LEU	2.7
2	B	475	GLY	2.7
1	C	216	THR	2.7
1	A	232	VAL	2.7
1	A	501	GLY	2.7
2	B	63	ASP	2.7
2	B	137	PHE	2.7
1	C	115	THR	2.7
2	B	324	THR	2.7
2	D	133	ARG	2.7
1	C	126	GLU	2.7
1	C	118	HIS	2.7
1	C	138	GLY	2.6
2	B	264	TRP	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	327	GLN	2.6
2	D	865	HIS	2.6
1	C	100	LEU	2.6
2	D	132	LYS	2.6
1	C	232	VAL	2.6
2	D	115	THR	2.6
1	A	118	HIS	2.5
1	A	779	HIS	2.5
1	C	117	HIS	2.5
1	A	471	PRO	2.5
1	A	631	LEU	2.5
1	A	168	ASN	2.5
2	D	50	ASN	2.5
1	A	268	GLU	2.5
1	C	50	ASN	2.5
1	A	230	SER	2.5
2	B	864	HIS	2.5
1	A	59	VAL	2.5
2	D	395	GLN	2.5
1	A	92	PRO	2.5
1	C	110	ASP	2.5
1	A	634	GLY	2.5
1	A	180	ILE	2.5
1	C	133	ARG	2.4
1	A	159	LEU	2.4
1	A	865	HIS	2.4
1	A	215	GLU	2.4
2	D	126	GLU	2.4
1	A	473	SER	2.4
1	A	91	ILE	2.4
1	C	534	ALA	2.4
1	C	127	LEU	2.4
1	C	229	LEU	2.4
2	D	228	GLY	2.4
1	A	102	LEU	2.4
2	B	134	LEU	2.4
2	B	123	ARG	2.4
1	A	476	GLU	2.4
2	B	78	PRO	2.4
1	C	865	HIS	2.4
1	A	172	VAL	2.4
1	C	233	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	62	THR	2.3
1	A	199	ALA	2.3
1	A	630	ALA	2.3
2	B	167	LYS	2.3
1	C	472	PHE	2.3
2	B	111	HIS	2.3
2	B	119	GLY	2.3
2	D	394	GLU	2.3
1	A	324	THR	2.3
1	A	864	HIS	2.3
1	A	93	LYS	2.3
1	C	178	LEU	2.3
2	B	62	THR	2.3
1	A	200	PHE	2.2
2	B	132	LYS	2.2
2	B	176	LYS	2.2
1	A	96	ARG	2.2
2	B	82	ARG	2.2
2	D	601	CYS	2.2
1	C	215	GLU	2.2
1	A	135	GLY	2.2
2	D	267	SER	2.2
2	B	601	CYS	2.2
1	A	397	LEU	2.2
1	A	197	ASP	2.2
2	B	474	PRO	2.2
2	D	598	ALA	2.2
1	C	864[A]	HIS	2.2
2	B	43	LYS	2.2
2	D	43	LYS	2.2
2	D	393	GLU	2.2
2	B	333	THR	2.2
2	D	230	SER	2.2
1	A	56	PHE	2.1
2	D	473	SER	2.1
2	D	166	GLY	2.1
2	B	100	LEU	2.1
2	B	174	PRO	2.1
2	B	397	LEU	2.1
1	C	176	LYS	2.1
2	D	130	GLU	2.1
2	B	64	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
2	B	116	PRO	2.1
2	B	128	LEU	2.1
2	B	135	GLY	2.1
1	A	632	GLN	2.1
1	C	434	GLU	2.0
1	A	600	SER	2.0
1	C	267	SER	2.0
1	A	136	VAL	2.0
2	B	20	PRO	2.0
1	A	512	LEU	2.0
1	C	123	ARG	2.0
1	C	129	ARG	2.0
1	A	88	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	CSO	A	844	7/8	0.91	0.09	35,38,47,52	0
1	CSO	C	844	7/8	0.94	0.08	34,40,52,53	0

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
4	EDO	D	914	4/4	0.79	0.22	54,54,57,63	0
4	EDO	A	907	4/4	0.80	0.17	55,55,55,57	0
4	EDO	D	913	4/4	0.82	0.15	49,56,57,58	0

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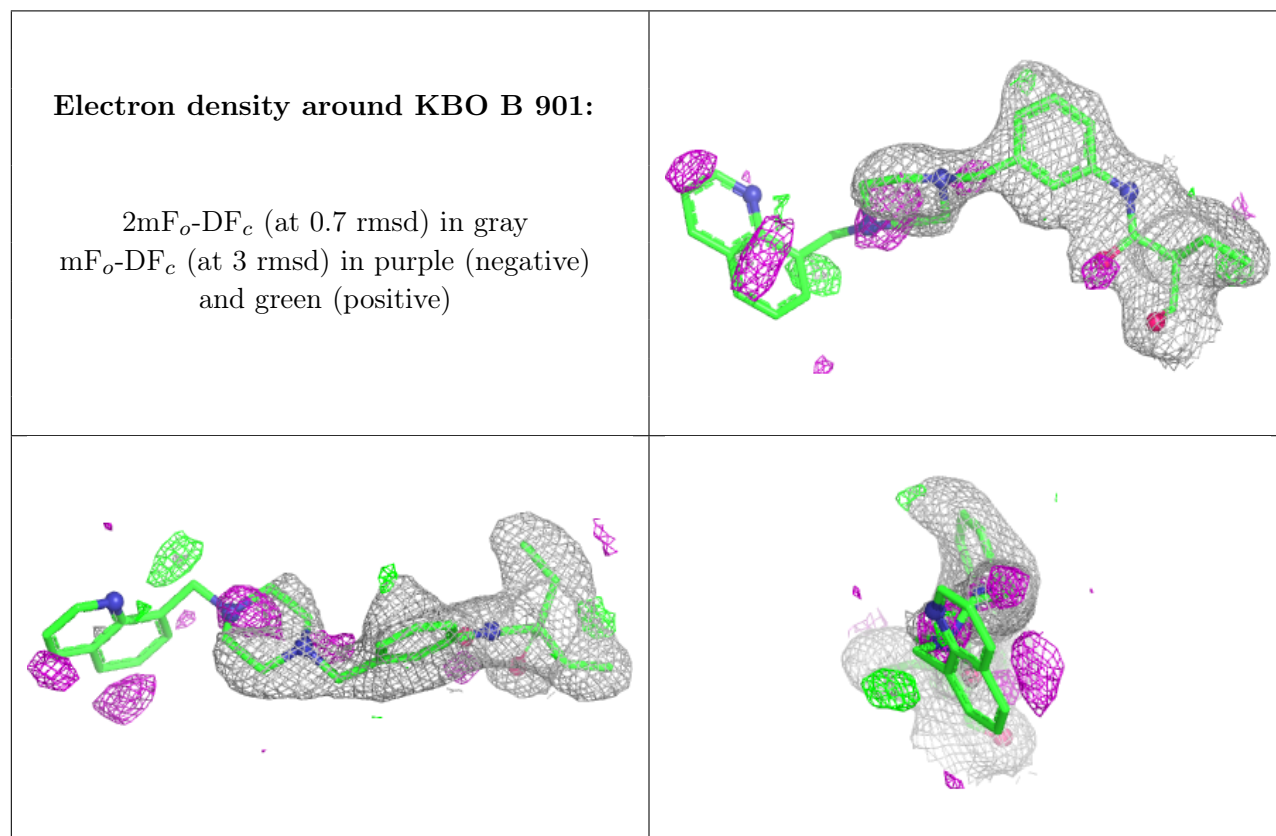
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	EDO	D	909	4/4	0.83	0.14	55,59,59,61	0
4	EDO	D	910	4/4	0.83	0.16	45,47,47,50	0
4	EDO	D	905	4/4	0.84	0.14	36,43,43,45	0
4	EDO	D	906	4/4	0.84	0.14	56,57,58,60	0
5	PEG	B	913	7/7	0.84	0.16	56,63,64,66	0
5	PEG	C	909	7/7	0.84	0.16	63,64,66,66	0
4	EDO	A	903	4/4	0.86	0.16	54,56,57,59	0
4	EDO	C	911	4/4	0.86	0.11	68,69,70,70	0
5	PEG	D	912	7/7	0.86	0.16	69,71,72,73	0
4	EDO	B	911	4/4	0.87	0.13	47,49,49,56	0
4	EDO	B	912	4/4	0.87	0.18	43,46,52,56	0
3	KBO	B	901	34/34	0.87	0.13	24,49,63,64	10
4	EDO	B	907	4/4	0.87	0.16	51,52,52,57	0
4	EDO	C	908	4/4	0.88	0.15	56,56,57,59	0
4	EDO	B	906	4/4	0.88	0.12	46,47,48,51	0
4	EDO	D	902	4/4	0.88	0.12	41,41,43,45	0
3	KBO	C	901	34/34	0.88	0.14	25,57,76,76	10
4	EDO	B	902	4/4	0.88	0.14	60,61,61,62	0
4	EDO	B	903	4/4	0.88	0.14	40,40,41,44	0
4	EDO	C	903	4/4	0.89	0.12	40,41,41,42	0
4	EDO	C	912	4/4	0.89	0.12	44,45,47,51	0
4	EDO	C	905	4/4	0.89	0.15	38,39,42,44	0
3	KBO	D	901	34/34	0.89	0.14	25,52,65,66	10
4	EDO	D	911	4/4	0.90	0.15	31,39,43,48	0
4	EDO	B	905	4/4	0.90	0.11	55,55,57,58	0
4	EDO	B	909	4/4	0.90	0.14	54,57,57,58	0
4	EDO	A	902	4/4	0.91	0.12	38,39,40,44	0
4	EDO	A	905	4/4	0.91	0.11	39,41,42,43	0
4	EDO	C	904	4/4	0.91	0.10	37,39,40,43	0
4	EDO	A	908	4/4	0.92	0.13	44,44,47,49	0
4	EDO	C	907	4/4	0.92	0.10	34,36,37,39	0
3	KBO	A	901	34/34	0.93	0.10	26,45,59,59	10
4	EDO	D	907	4/4	0.93	0.11	44,45,46,48	0
4	EDO	A	906	4/4	0.93	0.11	54,55,55,55	0
4	EDO	D	903	4/4	0.93	0.08	34,34,35,38	0
4	EDO	A	904	4/4	0.93	0.12	57,57,58,58	0
4	EDO	B	904	4/4	0.94	0.10	34,37,40,44	0
4	EDO	B	910	4/4	0.94	0.10	37,38,39,43	0
4	EDO	C	902	4/4	0.95	0.09	45,47,47,47	0
4	EDO	D	908	4/4	0.95	0.08	36,39,40,41	0
4	EDO	B	914	4/4	0.96	0.11	31,38,39,41	0
4	EDO	C	906	4/4	0.96	0.07	33,34,34,35	0

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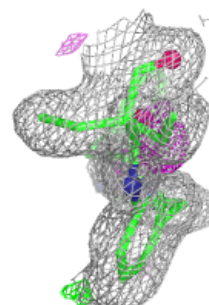
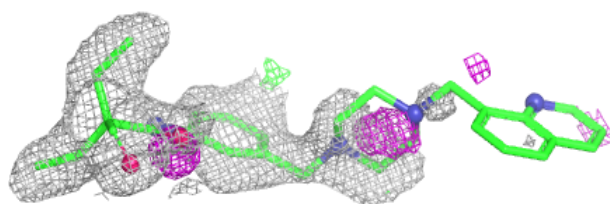
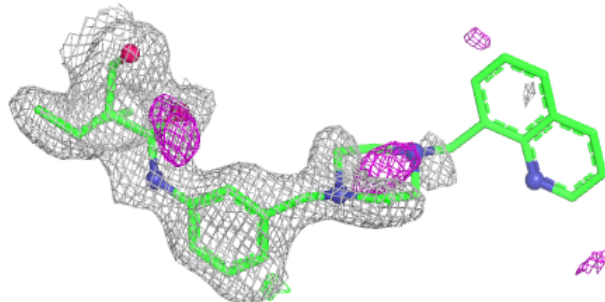
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	EDO	D	904	4/4	0.96	0.07	30,34,34,35	0
4	EDO	C	910	4/4	0.97	0.09	29,34,37,39	0
4	EDO	B	908	4/4	0.97	0.06	34,38,39,39	0

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.

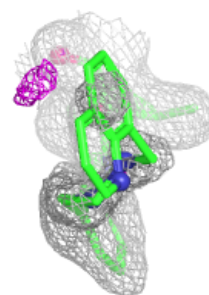
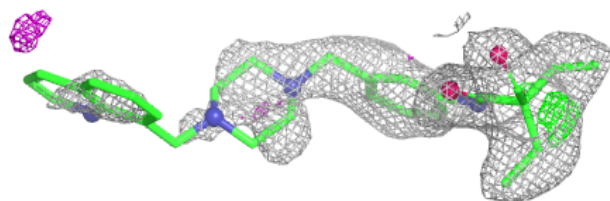
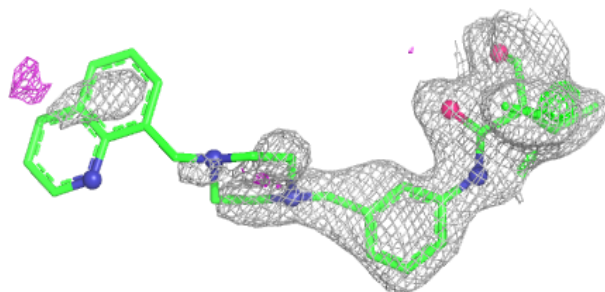


Electron density around KBO C 901:

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

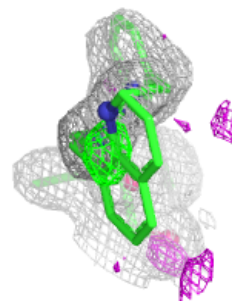
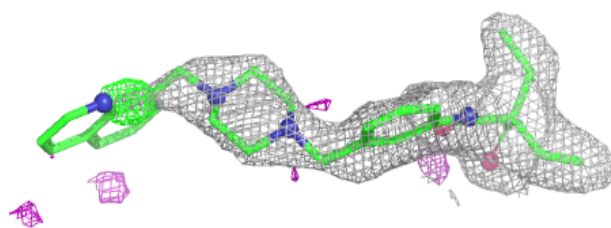
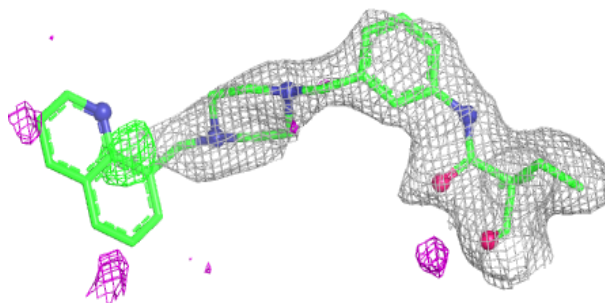
**Electron density around KBO D 901:**

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



Electron density around KBO A 901:

$2mF_o-DF_c$ (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.