



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:01 AM UTC

PDB ID : 6CWY / pdb_00006cwy
Title : Crystal structure of SUMO E1 in complex with an allosteric inhibitor
Authors : Lv, Z.; Yuan, L.; Atkison, J.H.; Williams, K.M.; Olsen, S.K.
Deposited on : 2018-04-01
Resolution : 2.46 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

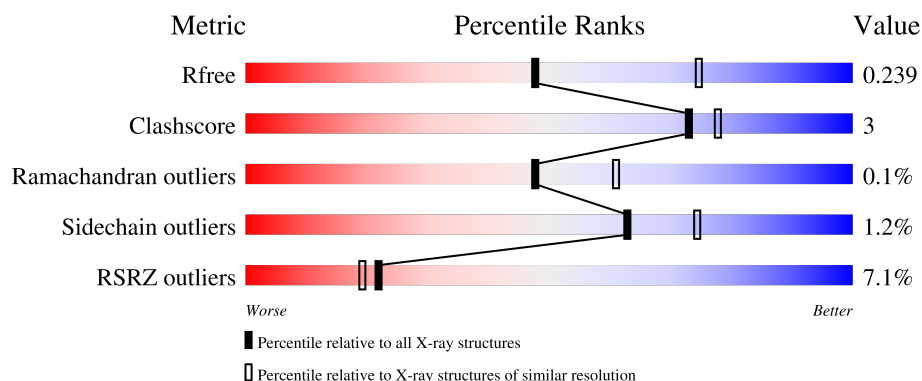
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

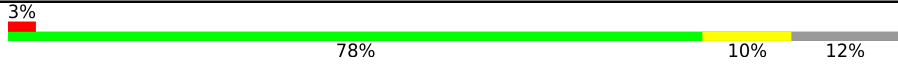
The reported resolution of this entry is 2.46 Å.

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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	C	346	
2	D	660	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6335 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 SUMO-activating enzyme subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	304	Total	C	N	O	S	0	1	0
			2402	1526	412	451	13			

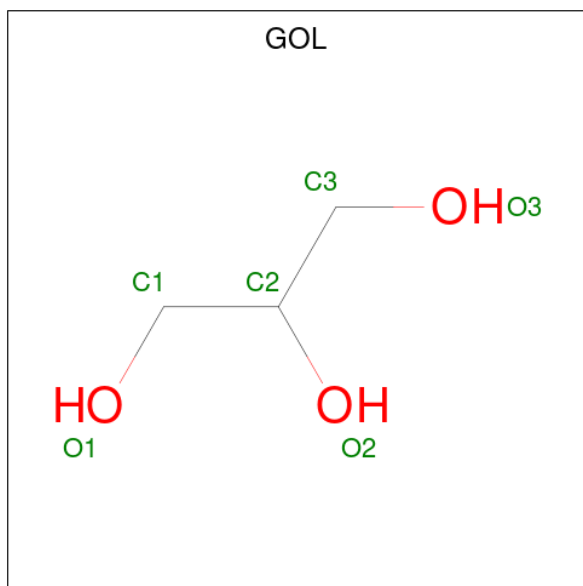
- Molecule 2 is a protein called SUMO-activating enzyme subunit 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	487	Total	C	N	O	S	0	1	0
			3823	2431	655	716	21			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-19	MET	-	initiating methionine	UNP Q9UBT2
D	-18	GLY	-	expression tag	UNP Q9UBT2
D	-17	SER	-	expression tag	UNP Q9UBT2
D	-16	SER	-	expression tag	UNP Q9UBT2
D	-15	HIS	-	expression tag	UNP Q9UBT2
D	-14	HIS	-	expression tag	UNP Q9UBT2
D	-13	HIS	-	expression tag	UNP Q9UBT2
D	-12	HIS	-	expression tag	UNP Q9UBT2
D	-11	HIS	-	expression tag	UNP Q9UBT2
D	-10	HIS	-	expression tag	UNP Q9UBT2
D	-9	SER	-	expression tag	UNP Q9UBT2
D	-8	SER	-	expression tag	UNP Q9UBT2
D	-7	GLY	-	expression tag	UNP Q9UBT2
D	-6	LEU	-	expression tag	UNP Q9UBT2
D	-5	VAL	-	expression tag	UNP Q9UBT2
D	-4	PRO	-	expression tag	UNP Q9UBT2
D	-3	ARG	-	expression tag	UNP Q9UBT2
D	-2	GLY	-	expression tag	UNP Q9UBT2
D	-1	SER	-	expression tag	UNP Q9UBT2
D	0	HIS	-	expression tag	UNP Q9UBT2

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

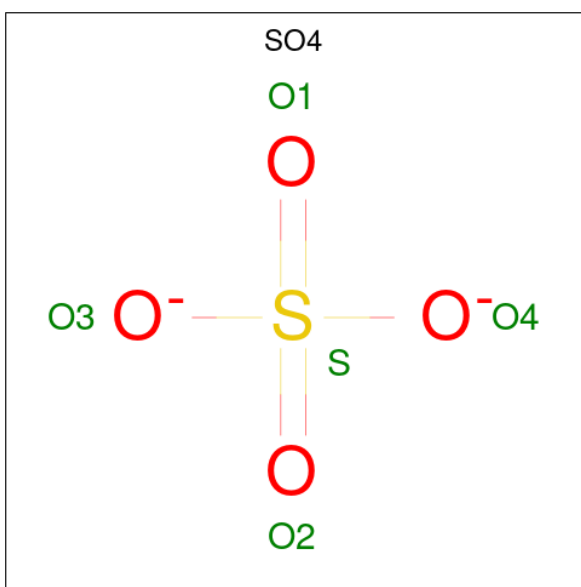
- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	1	Total	Zn	0	0
			1	1		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

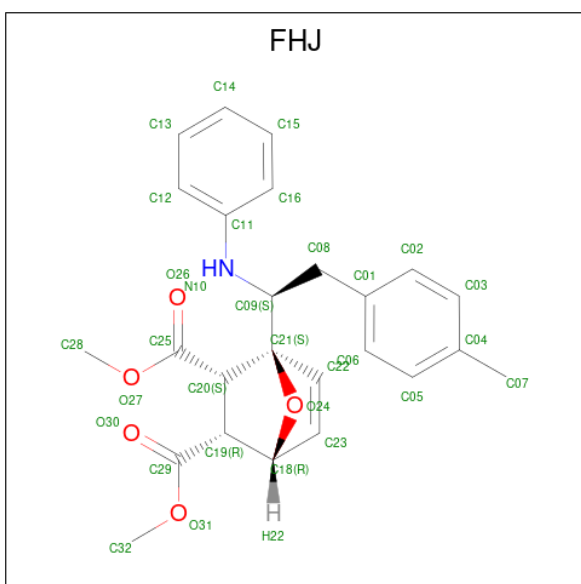
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	1	Total	Mg	0	0
			1	1		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	D	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is dimethyl (1S,2S,3R,4R)-1-[(1S)-2-(4-methylphenyl)-1-(phenylamino)ethyl]-7-oxabicyclo[2.2.1]hept-5-ene-2,3-dicarboxylate (CCD ID: FHJ) (formula: C₂₅H₂₇NO₅).

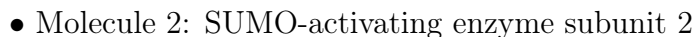


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	D	1	Total	C	N	O	0	0
			31	25	1	5		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	C	19	Total	O	0	0
			19	19		
8	D	26	Total	O	0	0
			26	26		

- Molecule 1: SUMO-activating enzyme subunit 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	56.05Å 116.05Å 174.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.28 – 2.46 48.28 – 2.46	Depositor EDS
% Data completeness (in resolution range)	99.3 (48.28-2.46) 94.4 (48.28-2.46)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 2.45Å)	Xtriage
Refinement program	PHENIX 1.10.1 _2155	Depositor
R, R_{free}	0.197 , 0.237 0.201 , 0.239	Depositor DCC
R_{free} test set	2000 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	48.3	Xtriage
Anisotropy	0.108	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 48.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6335	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.41% 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: FHJ, SO4, ZN, GOL, MG

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	C	0.08	0/2448	0.25	0/3300
2	D	0.09	0/3899	0.26	0/5280
All	All	0.08	0/6347	0.26	0/8580

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	C	2402	0	2412	17	0
2	D	3823	0	3854	23	0
3	C	12	0	16	1	0
4	D	1	0	0	0	0
5	D	1	0	0	0	0
6	D	20	0	0	0	0
7	D	31	0	0	0	0
8	C	19	0	0	0	0
8	D	26	0	0	0	0
All	All	6335	0	6282	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)
2:D:250:ASP:HB3	2:D:253:LYS:HB2	1.73	0.71
2:D:257:LYS:HA	2:D:261:ASP:HB2	1.75	0.69
2:D:256:THR:HG23	2:D:260:LYS:HD2	1.75	0.68
2:D:490:LEU:HD11	2:D:511:ILE:HD11	1.79	0.64
2:D:187:VAL:HG22	2:D:383:ILE:HD12	1.80	0.63
2:D:119:ARG:NH2	2:D:166:THR:OG1	2.35	0.60
2:D:363:MET:HE1	2:D:371:LYS:HG3	1.83	0.60
1:C:291:PHE:HA	3:C:401:GOL:H32	1.84	0.58
1:C:273:ARG:NH1	1:C:284:PRO:O	2.36	0.57
2:D:449:GLU:HG2	2:D:529:LEU:HB2	1.87	0.55
1:C:273:ARG:NH1	1:C:287:LEU:O	2.40	0.55
2:D:244:ALA:HB2	2:D:254:LEU:HD21	1.90	0.54
2:D:313:LEU:HD13	2:D:370:MET:HB2	1.90	0.54
2:D:537:ASP:O	2:D:539:GLY:N	2.40	0.54
2:D:462:LEU:HB2	2:D:502:ASN:HB3	1.91	0.53
1:C:224:TRP:CD1	1:C:282:ILE:HD11	2.45	0.52
1:C:217:LYS:O	1:C:221:GLU:HG2	2.10	0.52
2:D:52:ILE:HD13	2:D:70:ARG:HB3	1.92	0.51
2:D:52:ILE:HG12	2:D:93:HIS:HB2	1.93	0.51
2:D:260:LYS:HA	2:D:281:PRO:HB3	1.92	0.51
2:D:248:GLY:O	2:D:330:ARG:NH1	2.44	0.50
1:C:36:ALA:HA	1:C:61:LYS:HE3	1.92	0.50
1:C:120:GLU:HG3	1:C:145:ILE:HG23	1.96	0.48
1:C:163:GLY:HA3	1:C:306:VAL:HG21	1.94	0.47
2:D:348:ASP:HB3	2:D:351:ALA:HB3	1.97	0.47
1:C:30:ALA:HB1	1:C:319:SER:HB2	1.95	0.47
1:C:157:ASP:HB2	1:C:164:TYR:HE1	1.81	0.45
1:C:53:LYS:HE3	1:C:53:LYS:HB3	1.83	0.45
2:D:453:ARG:NH1	2:D:545:GLU:OE2	2.49	0.45
2:D:181:GLU:HG3	2:D:368:MET:HB2	2.00	0.44
2:D:256:THR:O	2:D:258:LEU:N	2.47	0.44
2:D:130:LEU:O	2:D:153:LYS:HD2	2.19	0.43
1:C:326:ASN:HA	1:C:327:ASN:HA	1.65	0.43
1:C:233:LEU:HD21	1:C:282:ILE:HG23	2.00	0.42
2:D:380:ALA:HA	2:D:383:ILE:HD11	2.02	0.42
2:D:137:ILE:HG21	2:D:398:ILE:HG23	2.02	0.42
1:C:259:SER:HG	1:C:297:SER:HG	1.63	0.41
1:C:160:GLY:HA2	1:C:299:MET:HE3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:161:TYR:CE1	1:C:299:MET:HE1	2.56	0.41
1:C:323:PRO:HA	1:C:324:PRO:HD3	1.94	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	C	301/346 (87%)	292 (97%)	9 (3%)	0	100	100
2	D	478/660 (72%)	457 (96%)	20 (4%)	1 (0%)	43	54
All	All	779/1006 (77%)	749 (96%)	29 (4%)	1 (0%)	48	61

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	538	LEU

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	C	264/296 (89%)	260 (98%)	4 (2%)	57	70
2	D	423/569 (74%)	418 (99%)	5 (1%)	63	74
All	All	687/865 (79%)	678 (99%)	9 (1%)	63	73

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	159	PHE
1	C	228	LYS
1	C	286	LEU
1	C	334	MET
2	D	30[A]	CYS
2	D	30[B]	CYS
2	D	247	THR
2	D	313	LEU
2	D	512	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	43	HIS
2	D	204	GLN

5.3.3 RNA [i](#)

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](#)

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 for Mogul analysis.

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	SO4	D	706	-	4,4,4	0.24	0	6,6,6	0.08	0
6	SO4	D	704	-	4,4,4	0.24	0	6,6,6	0.08	0
6	SO4	D	703	-	4,4,4	0.24	0	6,6,6	0.07	0
3	GOL	C	401	-	5,5,5	0.37	0	5,5,5	0.33	0
3	GOL	C	402	-	5,5,5	0.38	0	5,5,5	0.32	0
7	FHJ	D	707	2	33,34,34	1.14	2 (6%)	41,49,49	1.25	4 (9%)
6	SO4	D	705	-	4,4,4	0.23	0	6,6,6	0.08	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
3	GOL	C	401	-	-	2/4/4/4	-
3	GOL	C	402	-	-	2/4/4/4	-
7	FHJ	D	707	2	-	5/24/51/51	0/5/4/4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	D	707	FHJ	C11-N10	2.79	1.44	1.39
7	D	707	FHJ	C21-C22	2.03	1.54	1.51

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	707	FHJ	O31-C29-C19	3.72	116.33	111.03
7	D	707	FHJ	C20-C19-C29	3.45	121.07	114.00
7	D	707	FHJ	C20-C21-C22	2.33	111.47	107.61
7	D	707	FHJ	C21-C20-C19	-2.28	100.36	101.82

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	401	GOL	O1-C1-C2-C3
3	C	402	GOL	O1-C1-C2-C3
7	D	707	FHJ	C19-C29-O31-C32
7	D	707	FHJ	O30-C29-O31-C32

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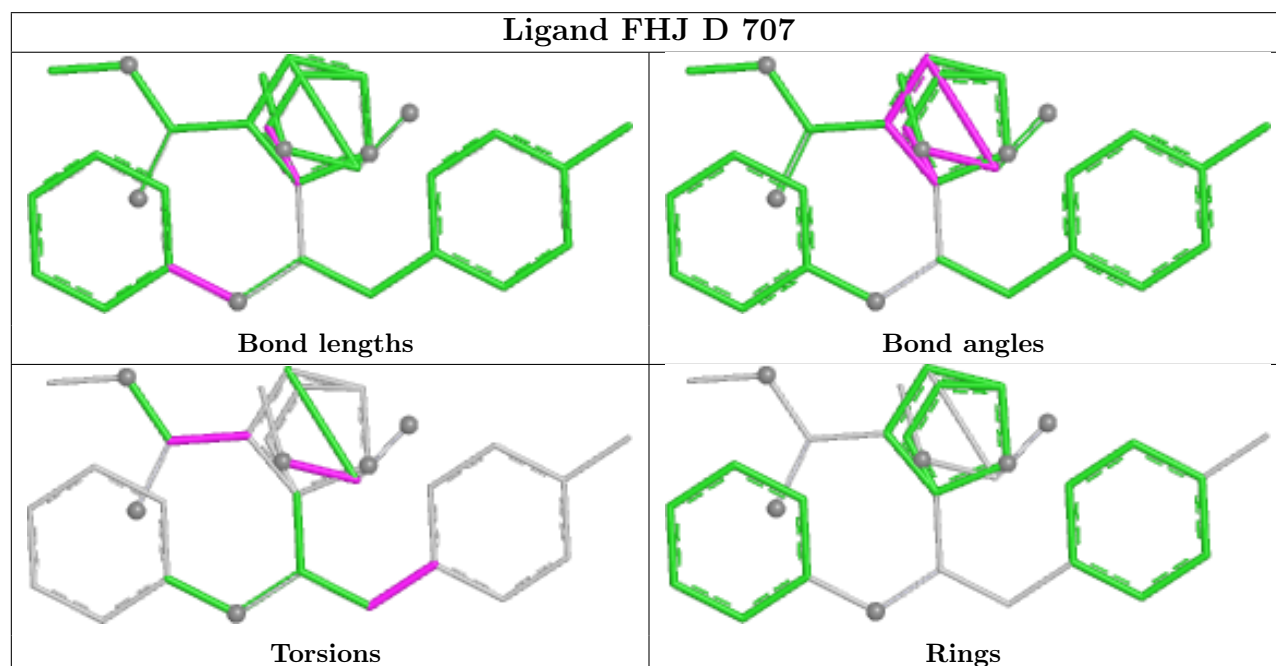
Mol	Chain	Res	Type	Atoms
3	C	401	GOL	O1-C1-C2-O2
3	C	402	GOL	O1-C1-C2-O2
7	D	707	FHJ	C06-C01-C08-C09
7	D	707	FHJ	C02-C01-C08-C09
7	D	707	FHJ	C19-C20-C25-O26

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	401	GOL	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	C	304/346 (87%)	0.22	9 (2%)	52	53	36, 63, 112, 152	1 (0%)
2	D	487/660 (73%)	0.47	47 (9%)	13	11	32, 63, 147, 199	1 (0%)
All	All	791/1006 (78%)	0.37	56 (7%)	22	19	32, 63, 140, 199	2 (0%)

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	432	CYS	6.4
2	D	69	GLY	6.2
2	D	224	ALA	6.1
2	D	238	ILE	6.0
2	D	315	VAL	5.5
2	D	380	ALA	4.7
2	D	342	LEU	4.3
1	C	205	THR	4.3
2	D	182	PRO	3.8
2	D	343	ILE	3.7
1	C	17	ALA	3.6
2	D	254	LEU	3.6
2	D	329	LEU	3.4
2	D	313	LEU	3.3
2	D	258	LEU	3.3
2	D	180	SER	3.2
1	C	230	LYS	3.1
2	D	197	PHE	3.1
2	D	346	LYS	3.1
1	C	346	LYS	3.0
2	D	344	TRP	2.9
2	D	257	LYS	2.9
1	C	178	LYS	2.8
2	D	240	THR	2.8

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Mol	Chain	Res	Type	RSRZ
2	D	282	LEU	2.7
2	D	368	MET	2.7
2	D	252	VAL	2.6
1	C	105	MET	2.6
2	D	322	PHE	2.6
1	C	286	LEU	2.5
2	D	379	MET	2.5
2	D	374	PHE	2.5
2	D	173	CYS	2.5
2	D	262	ASP	2.5
2	D	367	SER	2.5
2	D	548	GLY	2.4
2	D	196	LEU	2.4
2	D	50	ASP	2.4
2	D	216	ALA	2.4
2	D	165	PRO	2.3
2	D	284	TRP	2.3
2	D	287	VAL	2.3
2	D	275	LYS	2.3
2	D	52	ILE	2.2
1	C	107	ASP	2.2
2	D	274	ARG	2.2
2	D	255	PHE	2.2
2	D	331	VAL	2.2
2	D	280	VAL	2.2
2	D	381	GLY	2.2
2	D	212	ASP	2.1
2	D	421	GLN	2.1
2	D	260	LYS	2.1
2	D	281	PRO	2.1
1	C	25	LEU	2.0
2	D	181	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 [i](#)

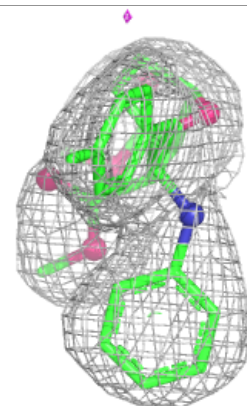
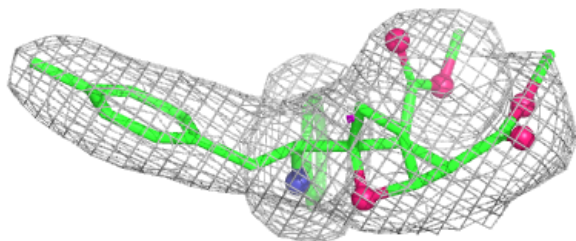
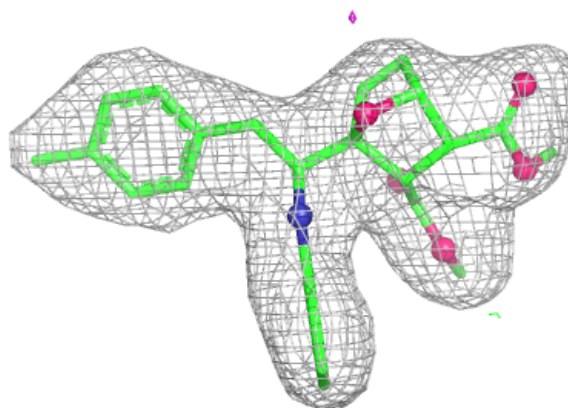
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	SO4	D	705	5/5	0.69	0.13	135,137,138,138	0
6	SO4	D	706	5/5	0.73	0.19	158,159,160,161	0
6	SO4	D	703	5/5	0.80	0.10	133,134,135,135	0
6	SO4	D	704	5/5	0.81	0.10	125,126,130,132	0
3	GOL	C	401	6/6	0.83	0.20	73,79,85,89	0
3	GOL	C	402	6/6	0.88	0.20	84,91,94,95	0
5	MG	D	702	1/1	0.95	0.22	60,60,60,60	0
7	FHJ	D	707	31/31	0.95	0.09	42,54,70,83	0
4	ZN	D	701	1/1	0.99	0.02	45,45,45,45	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 FHJ D 707:

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.