



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:08 AM UTC

PDB ID : 6JGR / pdb_00006jgr
Title : Crystal structure of barley exohydrolaseI W434Y in complex with 4'-nitrophenyl thiolaminaribioside
Authors : Luang, S.; Streltsov, V.A.; Hrmova, M.
Deposited on : 2019-02-14
Resolution : 2.25 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

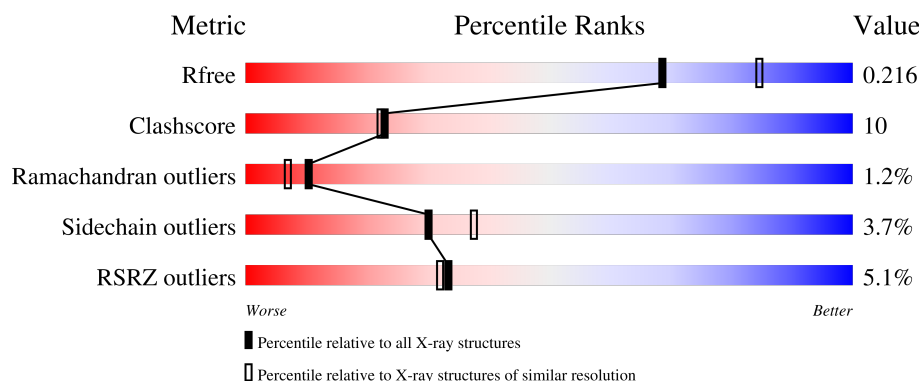
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	609	5% 83% 13% ...

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	1PE	A	705	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	ACT	A	724	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5076 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.

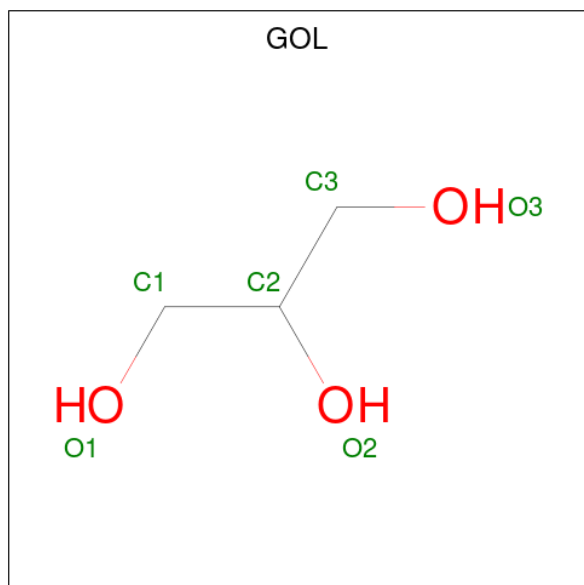
- Molecule 1 is a protein called BETA-D-GLUCAN GLUCOHYDROLASE ISOENZYME EXO1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	603	4591	2905	790	870	26	0	16	0

There are 6 discrepancies between the modelled and reference sequences:

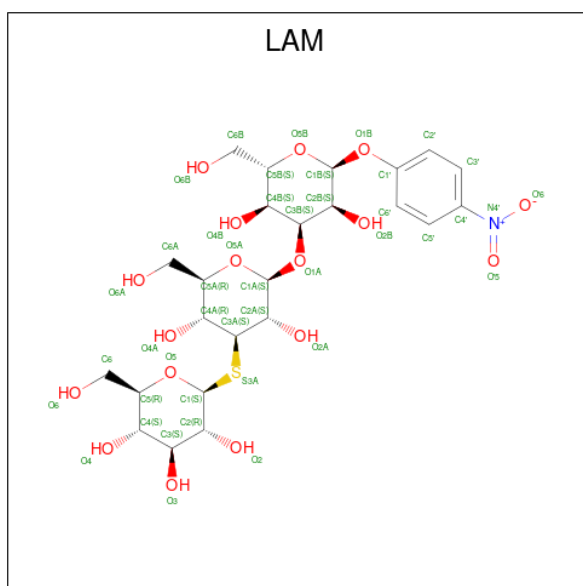
Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	HIS	-	expression tag	UNP A0A287SCR5
A	-2	HIS	-	expression tag	UNP A0A287SCR5
A	-1	ALA	-	expression tag	UNP A0A287SCR5
A	0	ALA	-	expression tag	UNP A0A287SCR5
A	320	LYS	ASN	See sequence details	UNP A0A287SCR5
A	434	TYR	TRP	engineered mutation	UNP A0A287SCR5

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



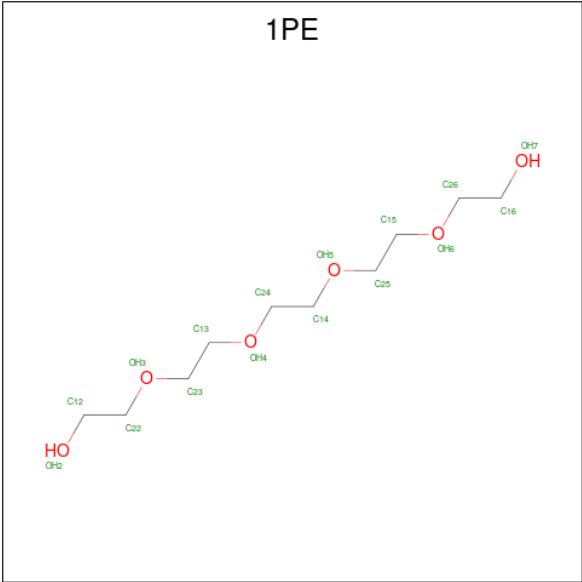
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is 4'-NITROPHENYL-S-(BETA-D-GLUCOPYRANOSYL)-(1-3)-(3-THIO-BETA-D-GLUCOPYRANOSYL)-(1-3)-BETA-D-GLUCOPYRANOSIDE (CCD ID: LAM) (formula: $C_{24}H_{35}NO_{17}S$).



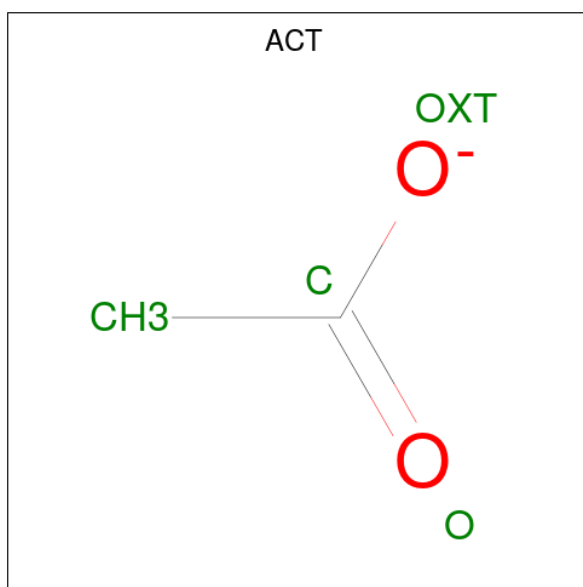
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	S	0	0
			23	12	10	1		

- Molecule 4 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			12	8	4		
4	A	1	Total	C	O	0	0
			9	6	3		
4	A	1	Total	C	O	0	0
			8	5	3		
4	A	1	Total	C	O	0	0
			9	6	3		
4	A	1	Total	C	O	0	0
			16	10	6		
4	A	1	Total	C	O	0	1
			12	7	5		
4	A	1	Total	C	O	0	0
			5	3	2		
4	A	1	Total	C	O	0	0
			12	8	4		

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



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

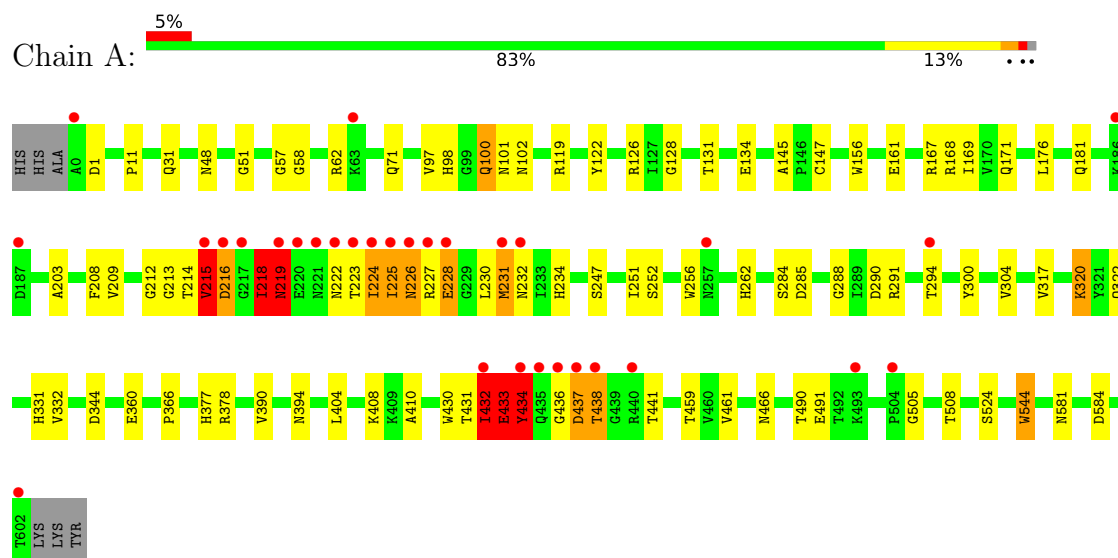
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	309	Total 309	O 309	0	0

3 Residue-property plots

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: BETA-D-GLUCAN GLUCOHYDROLASE ISOENZYME EXO1



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	100.47Å 100.47Å 179.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.85 – 2.25 45.85 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.8 (45.85-2.25) 100.0 (45.85-2.25)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.70 (at 2.24Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.163 , 0.208 0.174 , 0.216	Depositor DCC
R_{free} test set	2248 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	29.1	Xtriage
Anisotropy	0.142	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 57.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5076	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% 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: ACT, GOL, LAM, 1PE

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	1.26	13/4736 (0.3%)	1.17	15/6432 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	433	GLU	N-CA	7.00	1.55	1.46
1	A	317	VAL	CA-CB	6.18	1.57	1.54
1	A	432[A]	ILE	C-O	5.97	1.31	1.24
1	A	432[B]	ILE	C-O	5.97	1.31	1.24
1	A	544	TRP	CA-C	5.78	1.58	1.53
1	A	461	VAL	C-O	-5.73	1.18	1.24
1	A	219	ASN	CA-C	5.73	1.61	1.52
1	A	284	SER	C-O	5.47	1.31	1.23
1	A	432[A]	ILE	CA-C	5.38	1.59	1.52
1	A	432[B]	ILE	CA-C	5.38	1.59	1.52
1	A	128	GLY	N-CA	5.35	1.52	1.45
1	A	218	ILE	CA-C	5.19	1.59	1.52
1	A	390	VAL	N-CA	-5.17	1.40	1.46

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	219	ASN	N-CA-C	13.68	126.85	110.44
1	A	216	ASP	N-CA-C	11.54	126.86	111.28
1	A	433	GLU	N-CA-C	8.05	127.95	110.80
1	A	433	GLU	CA-C-N	6.01	132.52	121.70
1	A	433	GLU	C-N-CA	6.01	132.52	121.70
1	A	231	MET	N-CA-C	-5.88	106.19	113.19
1	A	436	GLY	N-CA-C	-5.77	104.35	112.60
1	A	434	TYR	CA-CB-CG	5.59	123.96	113.90
1	A	48	ASN	N-CA-C	5.48	119.66	112.92
1	A	432[A]	ILE	N-CA-C	5.43	120.64	109.34
1	A	432[B]	ILE	N-CA-C	5.43	120.64	109.34
1	A	332	VAL	N-CA-C	-5.39	105.59	110.82
1	A	360	GLU	N-CA-C	-5.19	107.08	113.41
1	A	438	THR	N-CA-C	-5.14	101.55	109.52
1	A	228	GLU	N-CA-C	-5.07	106.95	113.23

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	215	VAL	Peptide
1	A	219	ASN	Peptide
1	A	432[A]	ILE	Peptide
1	A	432[B]	ILE	Peptide

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	4591	0	4545	90	0
2	A	18	0	24	1	0
3	A	23	0	21	4	0
4	A	83	0	102	22	0
5	A	52	0	39	4	0
6	A	309	0	0	14	0
All	All	5076	0	4731	99	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 10.

All (99) 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:A:704:LAM:O5	6:A:805:HOH:O	1.80	0.98
1:A:97:VAL:HB	4:A:705:1PE:C23	1.94	0.98
1:A:433:GLU:HG2	1:A:434:TYR:HA	1.46	0.97
3:A:704:LAM:C1	6:A:805:HOH:O	2.12	0.97
1:A:466[A]:ASN:ND2	1:A:508:THR:OG1	1.98	0.95
1:A:394:ASN:HD21	1:A:404:LEU:H	1.16	0.93
1:A:433:GLU:CG	1:A:434:TYR:HA	1.99	0.92
3:A:704:LAM:S3A	6:A:805:HOH:O	2.31	0.88
1:A:219:ASN:HB2	1:A:491:GLU:HG3	1.56	0.86
1:A:97:VAL:H	1:A:101:ASN:HD21	1.35	0.74
1:A:145:ALA:HB1	4:A:705:1PE:C14	2.21	0.71
1:A:100:GLN:HE21	1:A:100:GLN:HA	1.55	0.70
1:A:218:ILE:HG22	1:A:218:ILE:O	1.92	0.68
1:A:408:LYS:H	5:A:717:ACT:H2	1.57	0.68
1:A:145:ALA:HB1	4:A:705:1PE:H142	1.75	0.67
4:A:709:1PE:H141	6:A:1030:HOH:O	1.96	0.65
1:A:156:TRP:HD1	1:A:219:ASN:HB3	1.62	0.64
1:A:100:GLN:HE21	1:A:100:GLN:CA	2.11	0.63
1:A:230:LEU:HB2	1:A:256:TRP:CZ2	2.33	0.63
4:A:709:1PE:C14	6:A:1030:HOH:O	2.46	0.63
1:A:11:PRO:HB3	4:A:711:1PE:H222	1.81	0.63
1:A:331:HIS:HD2	6:A:1062:HOH:O	1.81	0.62
1:A:212:GLY:O	1:A:218:ILE:HD13	1.99	0.62
4:A:709:1PE:C24	6:A:887:HOH:O	2.48	0.62
1:A:122:TYR:CE2	1:A:126:ARG:HD2	2.35	0.62
1:A:57:GLY:HA3	1:A:434:TYR:HB2	1.82	0.61
1:A:156:TRP:CD1	1:A:219:ASN:HB3	2.37	0.60
1:A:437:ASP:CB	1:A:441:THR:HG21	2.32	0.60
1:A:58:GLY:N	1:A:433:GLU:HB3	2.17	0.59
1:A:156:TRP:CG	1:A:218:ILE:HB	2.37	0.59
1:A:344:ASP:OD1	4:A:712:1PE:H121	2.01	0.59
1:A:214:THR:HB	1:A:215:VAL:HA	1.84	0.59
1:A:97:VAL:H	1:A:101:ASN:ND2	2.02	0.57
1:A:169:ILE:HG12	5:A:724:ACT:H1	1.85	0.57
1:A:262:HIS:HE1	1:A:285:ASP:H	1.51	0.57
1:A:134:GLU:OE2	1:A:377:HIS:HD2	1.88	0.56
1:A:145:ALA:HB1	4:A:705:1PE:H141	1.87	0.55
1:A:213:GLY:O	1:A:223:THR:HB	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:ASN:C	1:A:219:ASN:HD22	2.15	0.54
1:A:119:ARG:HD2	5:A:724:ACT:H2	1.88	0.54
1:A:97:VAL:HA	4:A:705:1PE:H252	1.91	0.53
1:A:181:GLN:HE21	1:A:203:ALA:H	1.54	0.53
1:A:58:GLY:H	1:A:102:ASN:ND2	2.08	0.52
1:A:234:HIS:HD2	6:A:972:HOH:O	1.90	0.52
1:A:432[B]:ILE:HD11	1:A:441:THR:HG21	1.91	0.52
1:A:169:ILE:HG12	5:A:724:ACT:CH3	2.41	0.51
1:A:300:TYR:CZ	1:A:320:LYS:HE2	2.46	0.51
1:A:490:THR:OG1	1:A:491:GLU:N	2.43	0.50
3:A:704:LAM:O4A	6:A:805:HOH:O	2.13	0.50
1:A:430:TRP:HE3	1:A:433:GLU:OE1	1.94	0.49
1:A:218:ILE:O	1:A:218:ILE:CG2	2.61	0.49
1:A:167:ARG:HH21	1:A:171:GLN:NE2	2.11	0.48
1:A:230:LEU:C	1:A:232:ASN:N	2.71	0.48
1:A:344:ASP:OD1	4:A:712:1PE:H222	2.13	0.48
1:A:122:TYR:CZ	1:A:126:ARG:HD2	2.47	0.48
1:A:581:ASN:HD22	4:A:708:1PE:H261	1.79	0.48
1:A:156:TRP:HB2	1:A:218:ILE:HB	1.96	0.47
4:A:707:1PE:C16	6:A:815:HOH:O	2.62	0.47
1:A:219:ASN:HB2	1:A:491:GLU:CG	2.37	0.47
4:A:709:1PE:H142	6:A:1030:HOH:O	2.12	0.47
1:A:100:GLN:HA	1:A:100:GLN:NE2	2.26	0.47
1:A:134:GLU:OE2	1:A:377:HIS:CD2	2.68	0.47
1:A:431:THR:C	1:A:432[B]:ILE:HG13	2.39	0.46
1:A:208:PHE:HA	1:A:209:VAL:HA	1.82	0.46
1:A:230:LEU:HD13	1:A:256:TRP:CE2	2.50	0.46
1:A:262:HIS:CE1	1:A:285:ASP:H	2.32	0.46
1:A:212:GLY:O	1:A:218:ILE:CD1	2.64	0.45
1:A:230:LEU:C	1:A:232:ASN:H	2.23	0.45
1:A:432[B]:ILE:HG12	1:A:441:THR:HB	1.98	0.45
1:A:71:GLN:HE22	1:A:366:PRO:HA	1.81	0.45
1:A:262:HIS:CE1	1:A:288:GLY:HA3	2.52	0.44
1:A:58:GLY:H	1:A:102:ASN:HD21	1.66	0.44
1:A:430:TRP:CZ2	1:A:434:TYR:CE1	3.06	0.44
1:A:98:HIS:HA	4:A:705:1PE:C16	2.48	0.43
1:A:234:HIS:CD2	6:A:972:HOH:O	2.67	0.43
1:A:433:GLU:CB	1:A:434:TYR:HA	2.45	0.43
1:A:145:ALA:CB	4:A:705:1PE:H142	2.44	0.43
1:A:31:GLN:HA	1:A:51:GLY:HA3	2.00	0.43
1:A:181:GLN:HE22	1:A:247:SER:H	1.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:THR:HG21	4:A:705:1PE:H152	2.01	0.42
1:A:524:SER:O	1:A:544:TRP:HA	2.20	0.42
1:A:176:LEU:HD21	4:A:705:1PE:H251	2.01	0.42
1:A:466[B]:ASN:OD1	1:A:508:THR:OG1	2.32	0.42
1:A:251:ILE:HG22	1:A:252:SER:N	2.34	0.42
1:A:161:GLU:OE2	4:A:705:1PE:C23	2.67	0.42
1:A:168:ARG:HA	1:A:171:GLN:HE21	1.84	0.42
1:A:394:ASN:HD22	1:A:394:ASN:HA	1.64	0.42
1:A:225:ILE:HG22	1:A:256:TRP:CZ2	2.54	0.42
1:A:432[A]:ILE:HD12	1:A:432[A]:ILE:HG23	1.61	0.41
1:A:300:TYR:O	1:A:304:VAL:HG23	2.20	0.41
1:A:147:CYS:HB3	4:A:705:1PE:H132	2.03	0.41
1:A:291:ARG:NH1	6:A:816:HOH:O	2.41	0.41
1:A:410:ALA:O	1:A:459:THR:HA	2.21	0.41
1:A:98:HIS:HA	4:A:705:1PE:H162	2.03	0.41
1:A:378:ARG:HD3	2:A:703:GOL:H11	2.02	0.41
1:A:230:LEU:HD23	1:A:231:MET:SD	2.61	0.40
1:A:320:LYS:HB2	6:A:1012:HOH:O	2.21	0.40
1:A:322[A]:GLN:OE1	4:A:709:1PE:H241	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	610/609 (100%)	576 (94%)	26 (4%)	8 (1%)	9 6

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	225	ILE
1	A	226	ASN

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Mol	Chain	Res	Type
1	A	433	GLU
1	A	432[A]	ILE
1	A	432[B]	ILE
1	A	505	GLY
1	A	224	ILE
1	A	218	ILE

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	494/490 (101%)	476 (96%)	18 (4%)	31	39

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

Mol	Chain	Res	Type
1	A	1	ASP
1	A	62	ARG
1	A	100	GLN
1	A	215	VAL
1	A	216	ASP
1	A	219	ASN
1	A	222	ASN
1	A	224	ILE
1	A	226	ASN
1	A	227	ARG
1	A	228	GLU
1	A	290	ASP
1	A	294	THR
1	A	320	LYS
1	A	433	GLU
1	A	434	TYR
1	A	437	ASP
1	A	438	THR

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	71	GLN
1	A	100	GLN
1	A	101	ASN
1	A	102	ASN
1	A	171	GLN
1	A	181	GLN
1	A	219	ASN
1	A	221	ASN
1	A	222	ASN
1	A	226	ASN
1	A	232	ASN
1	A	234	HIS
1	A	262	HIS
1	A	331	HIS
1	A	377	HIS
1	A	394	ASN
1	A	581	ASN

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](#)

26 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	$\# Z > 2$
5	ACT	A	716	-	3,3,3	0.92	0	3,3,3	1.04	0
4	1PE	A	705	-	11,11,15	0.61	0	10,10,14	1.16	0
5	ACT	A	715	-	3,3,3	0.62	0	3,3,3	1.24	0
5	ACT	A	724	-	3,3,3	0.82	0	3,3,3	2.60	2 (66%)
5	ACT	A	717	-	3,3,3	1.18	0	3,3,3	0.22	0
4	1PE	A	711	-	4,4,15	0.87	0	3,3,14	0.54	0
4	1PE	A	707	-	7,7,15	0.98	0	6,6,14	0.83	0
4	1PE	A	712	-	11,11,15	0.77	0	10,10,14	0.71	0
5	ACT	A	722	-	3,3,3	0.93	0	3,3,3	0.73	0
2	GOL	A	703	-	5,5,5	0.80	0	5,5,5	0.98	0
5	ACT	A	718	-	3,3,3	1.04	0	3,3,3	0.39	0
4	1PE	A	708	-	8,8,15	0.62	0	7,7,14	0.41	0
5	ACT	A	713	-	3,3,3	0.76	0	3,3,3	0.86	0
5	ACT	A	719	-	3,3,3	0.82	0	3,3,3	0.86	0
5	ACT	A	725	-	3,3,3	0.87	0	3,3,3	0.83	0
3	LAM	A	704	-	24,24,46	1.32	3 (12%)	27,35,67	1.85	5 (18%)
5	ACT	A	721	-	3,3,3	0.78	0	3,3,3	0.88	0
5	ACT	A	720	-	3,3,3	0.76	0	3,3,3	1.25	0
4	1PE	A	706	-	8,8,15	0.72	0	7,7,14	1.05	0
4	1PE	A	710[A]	-	7,7,15	0.62	0	6,6,14	0.61	0
2	GOL	A	702	-	5,5,5	0.50	0	5,5,5	1.15	0
5	ACT	A	723	-	3,3,3	1.09	0	3,3,3	1.14	0
5	ACT	A	714	-	3,3,3	0.98	0	3,3,3	0.48	0
2	GOL	A	701	-	5,5,5	0.63	0	5,5,5	1.50	1 (20%)
4	1PE	A	709	-	15,15,15	0.52	0	14,14,14	0.93	1 (7%)
4	1PE	A	710[B]	-	7,7,15	0.64	0	6,6,14	0.50	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	LAM	A	704	-	-	2/8/48/82	0/2/2/4
4	1PE	A	712	-	-	6/9/9/13	-
4	1PE	A	710[A]	-	-	3/5/5/13	-
2	GOL	A	702	-	-	2/4/4/4	-
2	GOL	A	703	-	-	1/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	1PE	A	708	-	-	3/6/6/13	-
4	1PE	A	709	-	-	8/13/13/13	-
4	1PE	A	710[B]	-	-	4/5/5/13	-
4	1PE	A	706	-	-	3/6/6/13	-
2	GOL	A	701	-	-	2/4/4/4	-
4	1PE	A	705	-	-	6/9/9/13	-
4	1PE	A	711	-	-	2/2/2/13	-
4	1PE	A	707	-	-	3/5/5/13	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	704	LAM	O5-C1	3.87	1.48	1.42
3	A	704	LAM	C4A-C3A	-2.21	1.51	1.53
3	A	704	LAM	C6-C5	2.11	1.58	1.51

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	704	LAM	C1-O5-C5	6.78	124.73	112.56
5	A	724	ACT	O-C-CH3	-3.43	108.48	122.53
3	A	704	LAM	O2-C2-C1	-2.86	105.22	110.30
5	A	724	ACT	OXT-C-CH3	2.83	126.93	115.05
4	A	709	1PE	OH7-C16-C26	-2.65	96.25	111.82
3	A	704	LAM	C4-C3-C2	2.17	114.64	110.83
2	A	701	GOL	O1-C1-C2	2.14	120.01	110.38
3	A	704	LAM	O5-C5-C6	2.05	111.51	106.44
3	A	704	LAM	C1-C2-C3	2.04	114.55	110.55

There are no chirality outliers.

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	702	GOL	O1-C1-C2-O2
2	A	702	GOL	O1-C1-C2-C3
4	A	710[A]	1PE	OH4-C13-C23-OH3
4	A	710[B]	1PE	OH4-C13-C23-OH3
4	A	708	1PE	OH6-C15-C25-OH5
4	A	709	1PE	OH2-C12-C22-OH3

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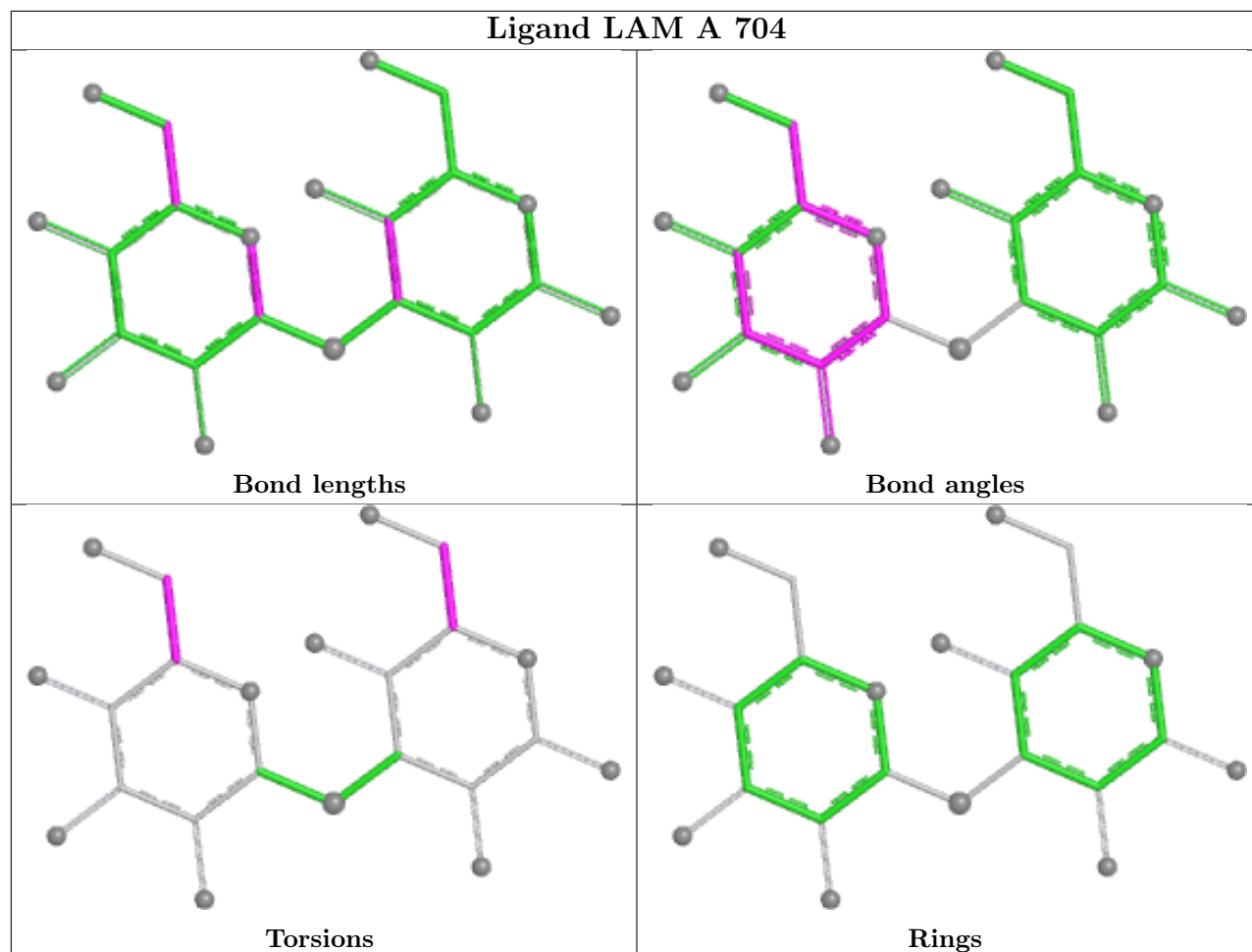
Mol	Chain	Res	Type	Atoms
4	A	705	1PE	OH5-C14-C24-OH4
4	A	709	1PE	OH5-C14-C24-OH4
4	A	712	1PE	OH5-C14-C24-OH4
4	A	707	1PE	C25-C15-OH6-C26
4	A	706	1PE	OH4-C13-C23-OH3
4	A	711	1PE	OH2-C12-C22-OH3
4	A	709	1PE	OH6-C15-C25-OH5
4	A	705	1PE	C25-C15-OH6-C26
2	A	703	GOL	C1-C2-C3-O3
4	A	705	1PE	OH7-C16-C26-OH6
4	A	712	1PE	OH2-C12-C22-OH3
4	A	707	1PE	OH6-C15-C25-OH5
3	A	704	LAM	O5A-C5A-C6A-O6A
4	A	709	1PE	OH4-C13-C23-OH3
4	A	705	1PE	OH6-C15-C25-OH5
4	A	706	1PE	OH2-C12-C22-OH3
4	A	705	1PE	C23-C13-OH4-C24
2	A	701	GOL	O2-C2-C3-O3
4	A	710[A]	1PE	C23-C13-OH4-C24
4	A	710[B]	1PE	C23-C13-OH4-C24
4	A	709	1PE	C14-C24-OH4-C13
4	A	710[B]	1PE	C12-C22-OH3-C23
4	A	712	1PE	C12-C22-OH3-C23
4	A	709	1PE	C23-C13-OH4-C24
4	A	708	1PE	C25-C15-OH6-C26
4	A	709	1PE	C24-C14-OH5-C25
4	A	710[A]	1PE	C12-C22-OH3-C23
4	A	709	1PE	OH7-C16-C26-OH6
4	A	710[B]	1PE	OH2-C12-C22-OH3
4	A	706	1PE	C14-C24-OH4-C13
4	A	711	1PE	C12-C22-OH3-C23
4	A	708	1PE	C24-C14-OH5-C25
4	A	707	1PE	OH7-C16-C26-OH6
4	A	712	1PE	C15-C25-OH5-C14
4	A	712	1PE	C24-C14-OH5-C25
2	A	701	GOL	O1-C1-C2-O2
4	A	705	1PE	C14-C24-OH4-C13
4	A	712	1PE	C14-C24-OH4-C13
3	A	704	LAM	O5-C5-C6-O6

There are no ring outliers.

10 monomers are involved in 31 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	705	1PE	12	0
5	A	724	ACT	3	0
5	A	717	ACT	1	0
4	A	711	1PE	1	0
4	A	707	1PE	1	0
4	A	712	1PE	2	0
2	A	703	GOL	1	0
4	A	708	1PE	1	0
3	A	704	LAM	4	0
4	A	709	1PE	5	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	603/609 (99%)	-0.11	31 (5%)	33 32	19, 28, 52, 85	20 (3%)

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	224	ILE	18.0
1	A	225	ILE	14.9
1	A	221	ASN	13.8
1	A	215	VAL	12.4
1	A	222	ASN	12.1
1	A	223	THR	10.4
1	A	217	GLY	9.4
1	A	227	ARG	9.1
1	A	226	ASN	8.8
1	A	220	GLU	8.4
1	A	228	GLU	6.7
1	A	0	ALA	6.5
1	A	216	ASP	6.3
1	A	434	TYR	4.9
1	A	435	GLN	3.7
1	A	231	MET	3.7
1	A	440	ARG	3.5
1	A	432[A]	ILE	3.2
1	A	438	THR	3.1
1	A	219	ASN	3.0
1	A	436	GLY	2.9
1	A	294	THR	2.9
1	A	187	ASP	2.8
1	A	437	ASP	2.5
1	A	186	LYS	2.5
1	A	602	THR	2.3
1	A	504	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	257	ASN	2.1
1	A	232	ASN	2.1
1	A	63	LYS	2.0
1	A	493	LYS	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.

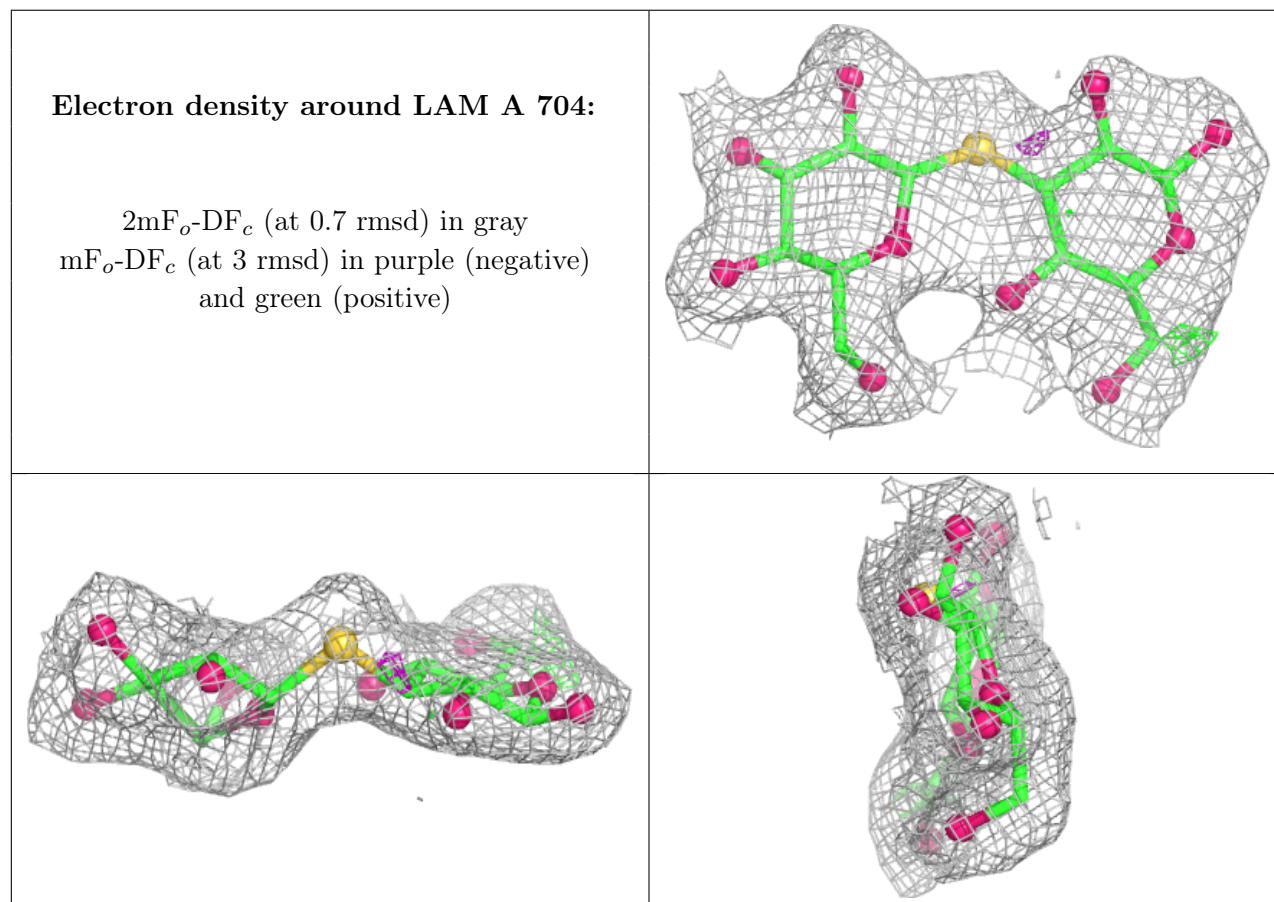
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	ACT	A	716	4/4	0.59	0.25	57,63,63,64	0
5	ACT	A	715	4/4	0.61	0.27	57,59,64,64	0
5	ACT	A	719	4/4	0.63	0.24	74,74,74,75	0
5	ACT	A	721	4/4	0.68	0.29	65,70,71,78	0
5	ACT	A	717	4/4	0.75	0.18	35,39,47,56	0
5	ACT	A	723	4/4	0.76	0.24	51,53,54,68	0
4	1PE	A	706	9/16	0.77	0.22	58,62,69,70	0
4	1PE	A	711	5/16	0.77	0.22	50,51,60,61	0
4	1PE	A	708	9/16	0.78	0.24	64,71,78,81	0
5	ACT	A	722	4/4	0.79	0.20	47,55,57,57	0
4	1PE	A	707	8/16	0.79	0.20	47,60,62,64	0
4	1PE	A	712	12/16	0.81	0.20	46,62,66,67	0
5	ACT	A	718	4/4	0.81	0.20	55,65,68,69	0
5	ACT	A	725	4/4	0.82	0.21	56,58,66,70	0
4	1PE	A	710[B]	8/16	0.85	0.20	39,57,67,68	4
4	1PE	A	710[A]	8/16	0.85	0.20	53,63,67,68	4
5	ACT	A	714	4/4	0.86	0.20	57,61,64,65	0
5	ACT	A	724	4/4	0.86	0.15	39,40,43,52	0
5	ACT	A	713	4/4	0.86	0.18	57,58,66,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	ACT	A	720	4/4	0.87	0.20	44,57,61,61	0
4	1PE	A	709	16/16	0.89	0.15	29,56,67,68	0
2	GOL	A	703	6/6	0.92	0.15	34,48,53,56	0
2	GOL	A	701	6/6	0.93	0.18	31,46,48,57	0
2	GOL	A	702	6/6	0.95	0.10	42,45,46,46	0
3	LAM	A	704	23/43	0.95	0.09	21,30,56,58	0
4	1PE	A	705	12/16	0.95	0.11	25,30,32,33	12

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.



6.5 Other polymers [i](#)

There are no such residues in this entry.