



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 13, 2023 – 11:23 am GMT

PDB ID : 4TPH
Title : Selectivity mechanism of a bacterial homologue of the human drug peptide transporters PepT1 and PepT2
Authors : Guettou, F.; Quistgaard, E.; Raba, M.; Moberg, P.; Low, C.; Nordlund, P.
Deposited on : 2014-06-07
Resolution : 3.15 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

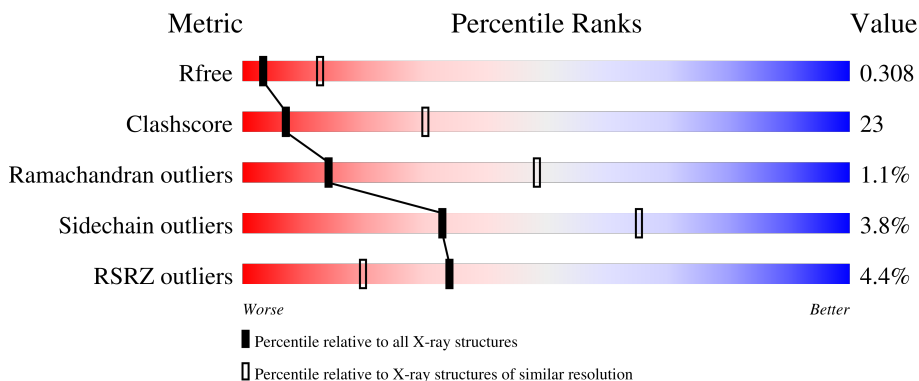
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.15 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}	130704	1665 (3.20-3.12)
Clashscore	141614	1804 (3.20-3.12)
Ramachandran outliers	138981	1770 (3.20-3.12)
Sidechain outliers	138945	1769 (3.20-3.12)
RSRZ outliers	127900	1616 (3.20-3.12)

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	523	
1	B	523	

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
3	LMT	B	602	-	-	-	X
4	ALA	A	603	-	-	-	X

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 6970 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 Proton:oligopeptide symporter POT family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	453	3422	2277	543	582	20	0	0	0
1	B	453	3422	2277	543	582	20	0	0	0

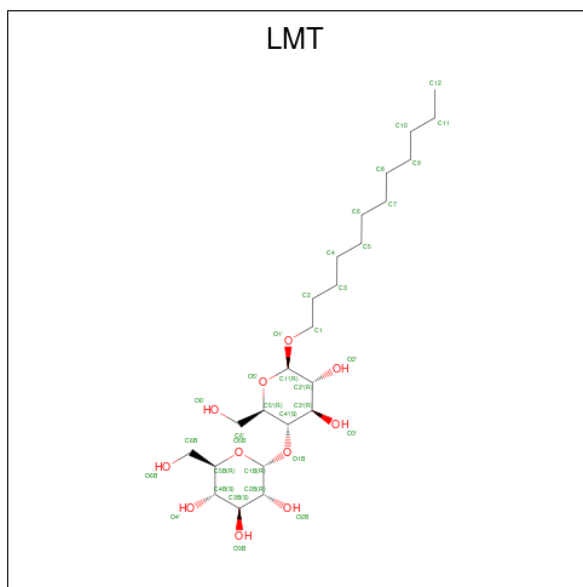
There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	517	ALA	-	expression tag	UNP Q8EHE6
A	518	GLU	-	expression tag	UNP Q8EHE6
A	519	ASN	-	expression tag	UNP Q8EHE6
A	520	LEU	-	expression tag	UNP Q8EHE6
A	521	TYR	-	expression tag	UNP Q8EHE6
A	522	PHE	-	expression tag	UNP Q8EHE6
A	523	GLN	-	expression tag	UNP Q8EHE6
B	517	ALA	-	expression tag	UNP Q8EHE6
B	518	GLU	-	expression tag	UNP Q8EHE6
B	519	ASN	-	expression tag	UNP Q8EHE6
B	520	LEU	-	expression tag	UNP Q8EHE6
B	521	TYR	-	expression tag	UNP Q8EHE6
B	522	PHE	-	expression tag	UNP Q8EHE6
B	523	GLN	-	expression tag	UNP Q8EHE6

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

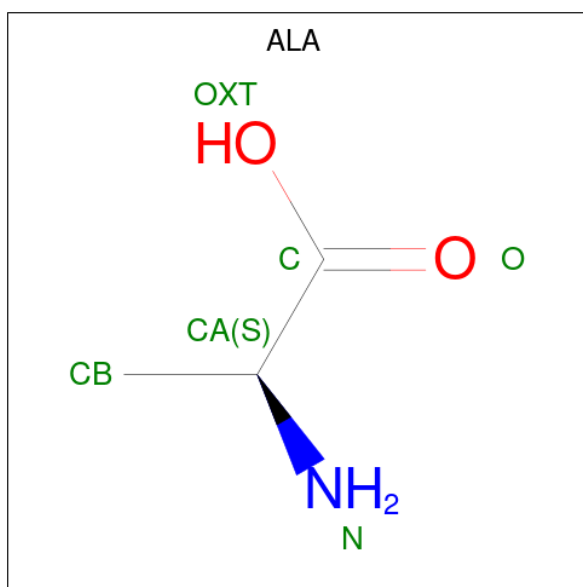
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is DODECYL-BETA-D-MALTOSE (three-letter code: LMT) (formula: C₂₄H₄₆O₁₁).



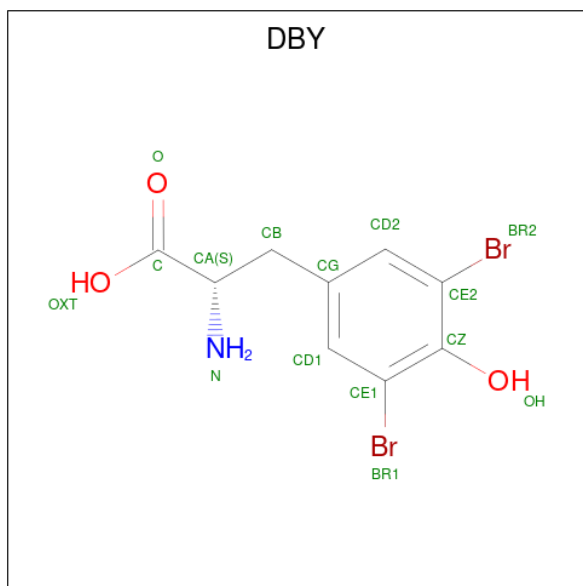
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			35	24	11		
3	B	1	Total	C	O	0	0
			35	24	11		
3	B	1	Total	C	O	0	0
			35	24	11		

- Molecule 4 is ALANINE (three-letter code: ALA) (formula: C₃H₇NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			5	3	1	1		

- Molecule 5 is 3,5 DIBROMOTYROSINE (three-letter code: DBY) (formula: $C_9H_9Br_2NO_3$).

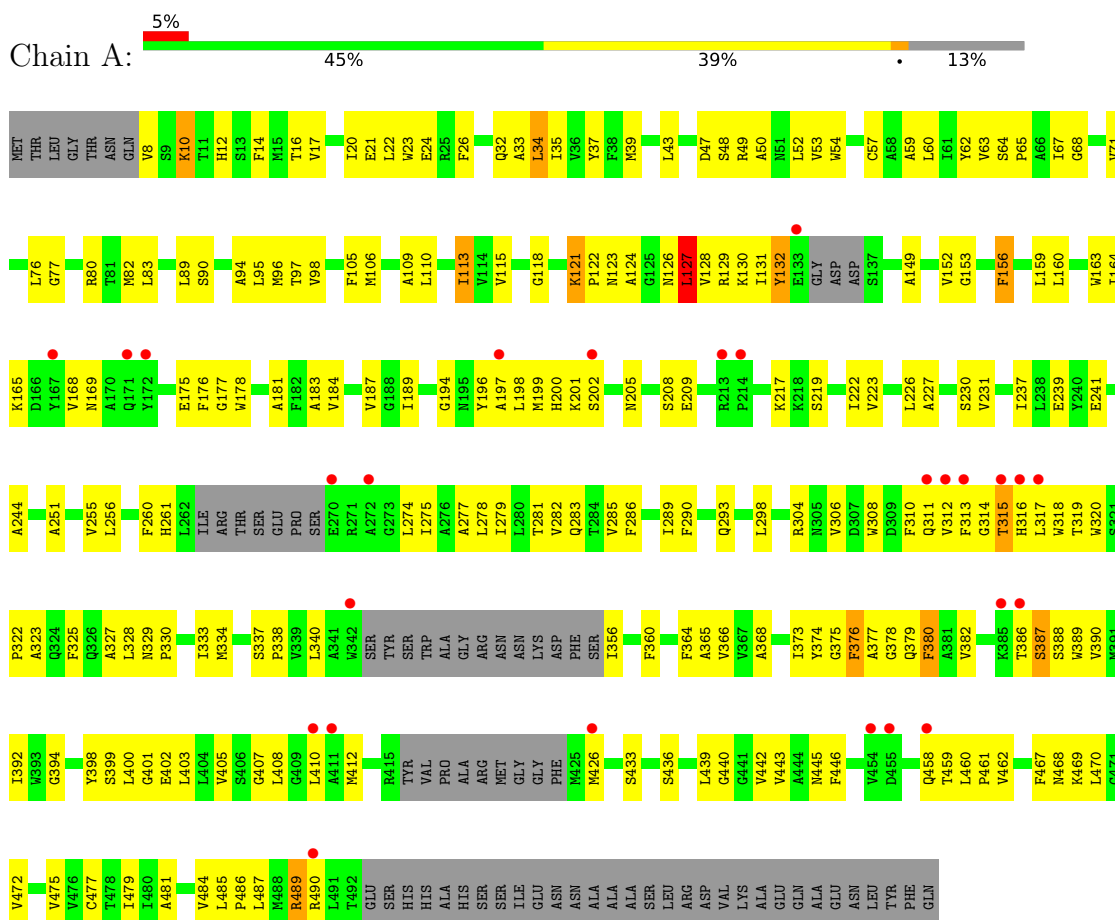


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	Br	C	N	O		
5	A	1	15	2	9	1	3	0	0

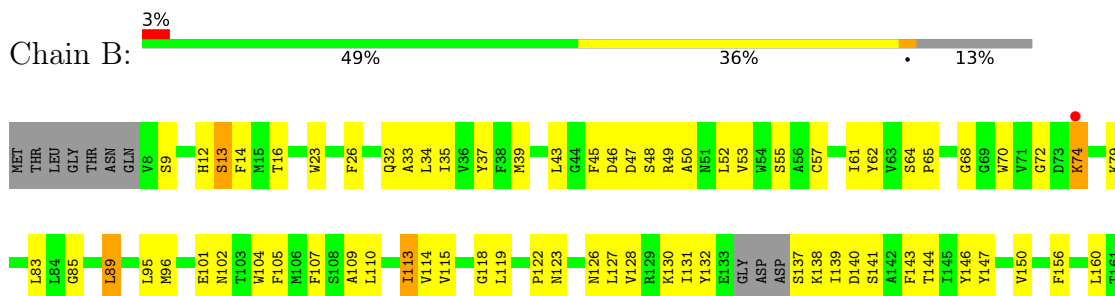
3 Residue-property plots i

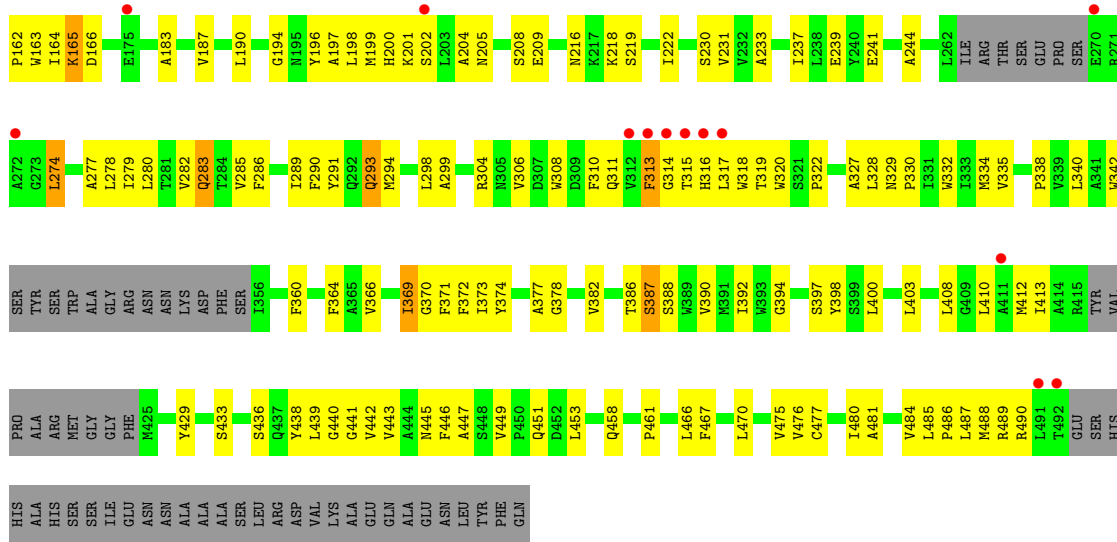
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Proton:oligopeptide symporter POT family



- Molecule 1: Proton:oligopeptide symporter POT family





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.65Å 107.68Å 204.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.54 – 3.15 29.54 – 3.16	Depositor EDS
% Data completeness (in resolution range)	78.2 (29.54-3.15) 78.4 (29.54-3.16)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 3.18Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.251 , 0.307 0.251 , 0.308	Depositor DCC
R_{free} test set	1256 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	101.4	Xtrriage
Anisotropy	0.059	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 70.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	6970	wwPDB-VP
Average B, all atoms (Å ²)	98.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% 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: ZN, DBY, LMT

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	0.64	1/3509 (0.0%)	0.85	4/4785 (0.1%)
1	B	0.63	1/3509 (0.0%)	0.82	3/4785 (0.1%)
All	All	0.64	2/7018 (0.0%)	0.83	7/9570 (0.1%)

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	1
1	B	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	33	ALA	C-O	5.64	1.34	1.23
1	B	33	ALA	C-O	5.27	1.33	1.23

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	127	LEU	CA-CB-CG	7.75	133.12	115.30
1	A	408	LEU	CA-CB-CG	6.61	130.50	115.30
1	B	408	LEU	CA-CB-CG	6.37	129.96	115.30
1	A	127	LEU	CB-CG-CD2	-5.89	100.99	111.00
1	A	34	LEU	CA-CB-CG	-5.38	102.94	115.30
1	B	89	LEU	CA-CB-CG	5.37	127.65	115.30
1	B	293	GLN	C-N-CA	-5.04	109.11	121.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	132	TYR	Peptide
1	B	274	LEU	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3422	0	3412	172	0
1	B	3422	0	3412	143	0
2	A	1	0	0	0	0
3	A	35	0	46	9	0
3	B	70	0	92	6	0
4	A	5	0	4	0	0
5	A	15	0	6	1	0
All	All	6970	0	6972	317	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 23.

All (317) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:364:PHE:CD2	1:B:477:CYS:HB3	2.03	0.93
1:A:152:VAL:HG22	1:A:156:PHE:CE1	2.08	0.89
1:A:94:ALA:O	1:A:97:THR:HG22	1.76	0.86
1:A:443:VAL:HA	1:A:446:PHE:HD2	1.42	0.84
1:B:200:HIS:CG	1:B:201:LYS:N	2.51	0.77
1:A:32:GLN:O	1:A:35:ILE:HG22	1.86	0.76
1:B:366:VAL:HG11	1:B:400:LEU:HD23	1.66	0.75
1:A:200:HIS:CG	1:A:201:LYS:N	2.55	0.75
1:A:76:LEU:HD11	1:A:223:VAL:HG21	1.68	0.74
1:A:219:SER:HA	1:A:222:ILE:HG22	1.69	0.74
1:A:20:ILE:HD12	1:A:196:TYR:HB2	1.70	0.74
1:A:382:VAL:HA	1:A:386:THR:HB	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:410:LEU:HD21	1:B:429:TYR:CD2	2.24	0.73
1:B:39:MET:HE1	1:B:53:VAL:HB	1.73	0.71
1:B:317:LEU:HD12	1:B:318:TRP:HA	1.73	0.71
1:A:156:PHE:HD1	3:A:602:LMT:H101	1.57	0.70
1:A:376:PHE:HA	1:A:379:GLN:OE1	1.91	0.70
1:A:374:TYR:HE2	1:A:398:TYR:CE2	2.10	0.69
1:B:197:ALA:O	1:B:200:HIS:ND1	2.24	0.69
1:B:477:CYS:HA	1:B:480:ILE:HD12	1.73	0.69
1:A:22:LEU:HD21	1:A:156:PHE:HE2	1.59	0.68
1:B:32:GLN:O	1:B:35:ILE:HG22	1.94	0.68
1:A:159:LEU:HB3	3:A:602:LMT:H52	1.74	0.68
1:B:13:SER:HA	1:B:16:THR:HG22	1.75	0.68
1:B:85:GLY:O	1:B:89:LEU:HD12	1.94	0.68
1:B:289:ILE:HG23	1:B:440:GLY:HA2	1.73	0.68
1:A:286:PHE:HA	1:A:289:ILE:HD12	1.76	0.68
1:B:382:VAL:HA	1:B:386:THR:HB	1.77	0.67
1:A:310:PHE:HB2	1:A:320:TRP:NE1	2.10	0.67
1:A:76:LEU:HB3	1:A:80:ARG:NH1	2.10	0.66
1:B:377:ALA:HB2	1:B:390:VAL:HG11	1.77	0.66
1:B:449:VAL:HG13	1:B:453:LEU:HD23	1.78	0.66
1:A:199:MET:HA	1:A:202:SER:CB	2.26	0.66
1:A:16:THR:HG22	1:A:199:MET:SD	2.36	0.66
1:B:488:MET:O	1:B:489:ARG:HD3	1.97	0.65
1:B:364:PHE:CE2	1:B:477:CYS:HB3	2.31	0.65
1:A:163:TRP:HE3	1:A:164:ILE:HD12	1.61	0.65
1:B:443:VAL:HA	1:B:446:PHE:HD2	1.62	0.65
1:A:315:THR:OG1	1:A:316:HIS:N	2.30	0.64
1:A:49:ARG:NH2	1:A:239:GLU:OE1	2.30	0.64
1:A:329:ASN:HB3	1:A:330:PRO:HD3	1.80	0.64
1:A:152:VAL:HG22	1:A:156:PHE:HE1	1.59	0.64
1:B:200:HIS:CG	1:B:201:LYS:H	2.14	0.64
1:A:156:PHE:CD1	3:A:602:LMT:H101	2.33	0.63
1:A:289:ILE:HG23	1:A:440:GLY:HA2	1.79	0.63
1:A:327:ALA:O	1:A:330:PRO:HD2	1.99	0.62
1:A:95:LEU:O	1:A:98:VAL:HG12	1.99	0.62
1:B:291:TYR:HA	1:B:398:TYR:OH	1.98	0.62
1:A:83:LEU:HD21	1:A:196:TYR:CD2	2.35	0.62
1:A:364:PHE:CD1	1:A:477:CYS:HB3	2.35	0.61
1:B:55:SER:HB3	1:B:438:TYR:HA	1.82	0.61
1:B:219:SER:HA	1:B:222:ILE:HG22	1.83	0.60
1:A:377:ALA:HB2	1:A:390:VAL:HG11	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:GLN:NE2	1:A:481:ALA:HB2	2.16	0.60
1:A:14:PHE:HB2	1:A:132:TYR:HE2	1.66	0.60
1:A:306:VAL:HG13	1:A:388:SER:H	1.66	0.60
1:B:308:TRP:O	1:B:320:TRP:N	2.35	0.60
1:B:306:VAL:HG13	1:B:388:SER:H	1.65	0.60
1:A:443:VAL:HA	1:A:446:PHE:CD2	2.30	0.59
1:A:197:ALA:O	1:A:200:HIS:N	2.32	0.59
1:A:365:ALA:O	1:A:368:ALA:HB3	2.02	0.59
1:A:373:ILE:HG22	1:A:390:VAL:HG13	1.83	0.59
1:B:34:LEU:HD11	1:B:162:PRO:HA	1.83	0.59
1:B:107:PHE:CD2	1:B:239:GLU:HG2	2.37	0.59
1:A:310:PHE:HB2	1:A:320:TRP:HE1	1.67	0.59
1:A:333:ILE:HG12	1:A:403:LEU:HD13	1.85	0.59
1:B:83:LEU:HD11	1:B:196:TYR:CE2	2.37	0.59
1:A:152:VAL:HG22	1:A:156:PHE:CZ	2.37	0.59
1:B:327:ALA:O	1:B:330:PRO:HD2	2.03	0.59
1:B:68:GLY:HA3	1:B:118:GLY:O	2.03	0.58
1:B:187:VAL:HA	1:B:190:LEU:HD12	1.83	0.58
1:B:39:MET:HE2	1:B:50:ALA:HA	1.83	0.58
1:A:96:MET:HG2	1:A:109:ALA:HB1	1.85	0.58
1:B:283:GLN:CD	1:B:360:PHE:HD2	2.06	0.58
1:B:147:TYR:O	1:B:150:VAL:HB	2.04	0.58
1:A:153:GLY:HA2	1:A:156:PHE:CE2	2.38	0.58
1:A:197:ALA:O	1:A:200:HIS:ND1	2.37	0.58
1:A:375:GLY:HA3	1:A:468:ASN:OD1	2.04	0.57
1:A:386:THR:O	1:A:387:SER:OG	2.20	0.57
1:A:65:PRO:HB3	1:A:122:PRO:HD3	1.86	0.57
1:B:328:LEU:HD13	1:B:392:ILE:HA	1.86	0.57
1:A:68:GLY:HA3	1:A:118:GLY:O	2.05	0.57
1:A:278:LEU:O	1:A:282:VAL:HG13	2.05	0.57
1:B:315:THR:CG2	1:B:316:HIS:H	2.17	0.57
1:A:39:MET:HE1	1:A:53:VAL:HB	1.86	0.57
1:A:48:SER:HA	1:A:445:ASN:OD1	2.04	0.57
1:A:67:ILE:O	1:A:71:VAL:HG23	2.04	0.57
1:B:48:SER:HA	1:B:445:ASN:OD1	2.05	0.56
1:B:274:LEU:HA	1:B:277:ALA:H	1.70	0.56
1:B:458:GLN:O	1:B:461:PRO:HD2	2.05	0.56
1:A:298:LEU:HD23	1:A:467:PHE:CZ	2.41	0.55
1:A:311:GLN:NE2	1:A:315:THR:H	2.04	0.55
1:A:311:GLN:NE2	1:A:315:THR:N	2.54	0.55
1:B:46:ASP:HB3	1:B:451:GLN:HE21	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:316:HIS:C	1:B:316:HIS:HD1	2.09	0.55
1:B:334:MET:O	1:B:338:PRO:HD3	2.06	0.55
1:A:479:ILE:HD12	1:B:369:ILE:HD12	1.88	0.55
1:B:298:LEU:HD23	1:B:467:PHE:CZ	2.42	0.55
1:A:181:ALA:O	1:A:184:VAL:HB	2.07	0.55
1:A:387:SER:HB3	1:A:390:VAL:HB	1.89	0.55
1:A:200:HIS:CG	1:A:201:LYS:H	2.25	0.55
1:A:256:LEU:HD22	1:A:260:PHE:HE2	1.72	0.54
1:B:138:LYS:HD2	1:B:141:SER:HB3	1.88	0.54
1:B:37:TYR:OH	1:B:299:ALA:O	2.20	0.54
1:A:128:VAL:HA	1:A:131:ILE:HD12	1.90	0.54
1:A:340:LEU:HD22	1:A:403:LEU:HD23	1.90	0.54
1:B:197:ALA:O	1:B:200:HIS:N	2.33	0.54
1:A:374:TYR:CE2	1:A:398:TYR:CE2	2.93	0.53
1:A:308:TRP:O	1:A:320:TRP:N	2.36	0.53
1:B:329:ASN:HB3	1:B:330:PRO:HD3	1.91	0.53
1:B:237:ILE:O	1:B:244:ALA:HB2	2.07	0.53
1:A:160:LEU:HG	1:A:164:ILE:HD13	1.90	0.53
1:A:366:VAL:HG11	1:A:400:LEU:HD23	1.91	0.53
1:A:183:ALA:O	1:A:187:VAL:HG13	2.08	0.53
1:B:447:ALA:HB2	1:B:466:LEU:HB2	1.91	0.53
1:B:137:SER:OG	1:B:138:LYS:N	2.39	0.53
1:B:132:TYR:HB2	1:B:139:ILE:HG22	1.90	0.52
1:B:279:ILE:O	1:B:282:VAL:HG22	2.09	0.52
1:A:475:VAL:HG21	1:B:475:VAL:HG21	1.90	0.52
1:B:199:MET:HA	1:B:202:SER:CB	2.40	0.52
1:B:283:GLN:NE2	1:B:360:PHE:HD2	2.08	0.52
3:B:602:LMT:H1B	3:B:602:LMT:O3'	2.08	0.52
1:B:101:GLU:OE1	1:B:101:GLU:N	2.42	0.52
1:B:386:THR:O	1:B:387:SER:OG	2.21	0.52
1:B:286:PHE:HA	1:B:289:ILE:HD12	1.91	0.52
1:B:316:HIS:C	1:B:316:HIS:ND1	2.63	0.52
1:A:109:ALA:O	1:A:113:ILE:HD12	2.11	0.51
1:A:442:VAL:O	1:A:445:ASN:HB2	2.11	0.51
1:A:489:ARG:HG3	1:A:489:ARG:O	2.10	0.51
1:B:315:THR:HG23	1:B:316:HIS:H	1.74	0.51
1:A:379:GLN:H	1:A:379:GLN:CD	2.14	0.51
1:A:310:PHE:HB2	1:A:320:TRP:CD1	2.45	0.51
1:A:378:GLY:C	1:A:380:PHE:H	2.13	0.51
1:B:65:PRO:HB3	1:B:122:PRO:HD3	1.92	0.51
1:A:328:LEU:HD13	1:A:392:ILE:HA	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:VAL:HG11	1:A:105:PHE:CZ	2.46	0.50
1:A:115:VAL:HG21	1:A:231:VAL:HG23	1.93	0.50
1:B:79:LYS:HD3	1:B:204:ALA:HB1	1.93	0.50
1:B:143:PHE:O	1:B:146:TYR:HB3	2.10	0.50
1:A:459:THR:O	1:A:462:VAL:HG22	2.11	0.50
1:B:315:THR:CG2	1:B:316:HIS:N	2.73	0.50
1:B:9:SER:HB2	1:B:12:HIS:CD2	2.46	0.50
1:B:62:TYR:CE2	1:B:433:SER:HB2	2.47	0.49
1:A:39:MET:HE2	1:A:50:ALA:HA	1.94	0.49
1:A:460:LEU:HB3	1:A:461:PRO:HD3	1.94	0.49
1:B:16:THR:OG1	1:B:199:MET:HB3	2.12	0.49
1:B:410:LEU:HD21	1:B:429:TYR:CE2	2.47	0.49
1:B:481:ALA:O	1:B:484:VAL:HG12	2.12	0.49
1:B:410:LEU:N	1:B:410:LEU:HD23	2.27	0.49
1:B:485:LEU:N	1:B:486:PRO:HD2	2.27	0.49
1:A:76:LEU:HA	1:A:80:ARG:NH2	2.27	0.49
1:A:274:LEU:HA	1:A:277:ALA:H	1.77	0.49
1:A:379:GLN:OE1	1:A:379:GLN:N	2.34	0.49
1:A:127:LEU:O	1:A:131:ILE:HG13	2.13	0.49
1:B:290:PHE:O	1:B:293:GLN:HB2	2.12	0.49
1:A:21:GLU:OE2	1:A:121:LYS:HE2	2.12	0.49
1:A:237:ILE:HG23	1:A:244:ALA:HA	1.94	0.49
1:A:12:HIS:O	1:A:16:THR:HG23	2.13	0.49
1:A:251:ALA:O	1:A:255:VAL:HG13	2.13	0.48
1:B:412:MET:SD	1:B:412:MET:N	2.86	0.48
1:A:98:VAL:HG11	1:A:105:PHE:CE1	2.48	0.48
1:A:115:VAL:HG11	1:A:227:ALA:HB1	1.96	0.48
1:B:64:SER:HB3	1:B:65:PRO:HD3	1.95	0.48
1:B:47:ASP:OD2	1:B:304:ARG:NH1	2.47	0.48
1:A:319:THR:OG1	3:A:602:LMT:H3B	2.13	0.48
1:B:55:SER:OG	1:B:441:GLY:HA3	2.14	0.47
1:B:458:GLN:C	1:B:461:PRO:HD2	2.35	0.47
1:B:57:CYS:SG	1:B:110:LEU:HD22	2.54	0.47
1:B:197:ALA:HA	1:B:200:HIS:HB3	1.96	0.47
1:B:128:VAL:HA	1:B:131:ILE:HD12	1.96	0.47
1:A:469:LYS:HE2	1:A:469:LYS:HB3	1.44	0.47
1:B:319:THR:OG1	3:B:601:LMT:H3B	2.14	0.47
1:A:60:LEU:HD23	1:A:60:LEU:HA	1.52	0.47
1:A:285:VAL:HG13	1:A:436:SER:HB3	1.97	0.47
1:A:394:GLY:O	1:A:398:TYR:HD2	1.97	0.47
1:B:127:LEU:O	1:B:131:ILE:HG13	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:387:SER:HB3	1:B:390:VAL:HB	1.97	0.47
1:A:334:MET:O	1:A:338:PRO:HD3	2.15	0.47
1:A:374:TYR:HE2	1:A:398:TYR:CZ	2.31	0.47
1:B:374:TYR:CZ	1:B:394:GLY:HA3	2.50	0.47
1:A:163:TRP:CE3	1:A:164:ILE:HD12	2.47	0.46
1:A:277:ALA:O	1:A:281:THR:HG23	2.15	0.46
1:B:283:GLN:NE2	1:B:360:PHE:CD2	2.83	0.46
1:B:370:GLY:O	1:B:373:ILE:HB	2.15	0.46
1:A:23:TRP:O	1:A:26:PHE:HB3	2.16	0.46
1:A:153:GLY:HA2	1:A:156:PHE:CD2	2.51	0.46
1:A:165:LYS:O	1:A:169:ASN:HB2	2.16	0.46
1:B:102:ASN:HD21	3:B:602:LMT:H3'	1.79	0.46
1:A:159:LEU:HD23	1:A:159:LEU:HA	1.65	0.46
1:A:194:GLY:O	1:A:198:LEU:HG	2.15	0.46
1:B:104:TRP:CE2	1:B:239:GLU:HG3	2.50	0.46
1:B:14:PHE:HB2	1:B:132:TYR:HE2	1.81	0.46
1:B:313:PHE:HA	1:B:314:GLY:HA2	1.50	0.46
1:A:311:GLN:HE21	1:A:314:GLY:HA2	1.81	0.46
1:A:458:GLN:O	1:A:461:PRO:HD2	2.15	0.46
1:B:280:LEU:HD23	1:B:280:LEU:HA	1.78	0.46
1:A:62:TYR:HE2	1:A:433:SER:HB2	1.81	0.46
1:A:64:SER:OG	1:A:118:GLY:HA3	2.16	0.46
1:A:410:LEU:HD22	1:A:426:MET:HA	1.98	0.46
1:B:233:ALA:O	1:B:237:ILE:HG13	2.16	0.46
1:B:310:PHE:HB2	1:B:320:TRP:CD1	2.51	0.46
1:A:34:LEU:HD23	1:A:34:LEU:HA	1.74	0.45
1:A:39:MET:CE	1:A:50:ALA:HA	2.46	0.45
1:A:52:LEU:HG	1:A:445:ASN:HD21	1.80	0.45
1:B:194:GLY:O	1:B:198:LEU:HG	2.17	0.45
1:B:83:LEU:HD21	1:B:196:TYR:CD2	2.51	0.45
1:B:216:ASN:HD21	1:B:218:LYS:NZ	2.14	0.45
1:B:279:ILE:HG21	1:B:484:VAL:HG21	1.97	0.45
1:B:476:VAL:HG12	1:B:480:ILE:HD11	1.98	0.45
1:A:126:ASN:OD1	1:A:130:LYS:NZ	2.49	0.45
1:A:196:TYR:O	1:A:196:TYR:CG	2.69	0.45
1:B:183:ALA:O	1:B:187:VAL:HG13	2.17	0.45
1:B:278:LEU:O	1:B:282:VAL:HG13	2.15	0.45
1:A:168:VAL:HB	1:A:177:GLY:HA2	1.99	0.45
1:A:47:ASP:OD2	1:A:304:ARG:NH1	2.50	0.45
1:A:389:TRP:CE3	1:A:392:ILE:HD12	2.51	0.45
1:A:394:GLY:O	1:A:398:TYR:CD2	2.70	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:LEU:HA	1:B:318:TRP:HA	1.57	0.45
3:B:601:LMT:H2'	3:B:601:LMT:H12	1.53	0.45
1:A:275:ILE:O	1:A:279:ILE:HG12	2.17	0.45
1:B:294:MET:SD	1:B:398:TYR:CE2	3.10	0.45
1:B:311:GLN:HB2	1:B:315:THR:O	2.17	0.45
1:B:62:TYR:HE2	1:B:433:SER:HB2	1.81	0.44
1:A:126:ASN:O	1:A:130:LYS:HD3	2.17	0.44
1:B:52:LEU:HD23	1:B:52:LEU:HA	1.43	0.44
1:A:8:VAL:CG1	1:A:12:HIS:CD2	3.00	0.44
1:A:156:PHE:CD1	3:A:602:LMT:H121	2.53	0.44
1:A:77:GLY:H	1:A:80:ARG:CG	2.30	0.44
1:A:89:LEU:CD2	1:A:113:ILE:HA	2.48	0.44
1:B:310:PHE:HB2	1:B:320:TRP:NE1	2.32	0.44
1:A:124:ALA:O	1:A:128:VAL:HG23	2.17	0.44
1:B:340:LEU:HD22	1:B:403:LEU:HD23	1.99	0.44
1:A:57:CYS:SG	1:A:110:LEU:HD22	2.58	0.44
1:A:115:VAL:HG22	1:A:230:SER:HB2	1.99	0.44
1:A:322:PRO:HA	1:A:325:PHE:CD2	2.53	0.43
1:A:402:GLU:HG3	5:A:604:DBY:BR2	2.73	0.43
1:A:35:ILE:HG21	1:A:54:TRP:CE3	2.53	0.43
1:B:487:LEU:HD23	1:B:487:LEU:HA	1.92	0.43
1:A:281:THR:O	1:A:285:VAL:HG23	2.18	0.43
1:B:39:MET:O	1:B:45:PHE:HB2	2.18	0.43
1:B:160:LEU:HG	1:B:164:ILE:HD13	1.99	0.43
1:B:109:ALA:O	1:B:113:ILE:HD12	2.17	0.43
1:B:342:TRP:CD1	1:B:342:TRP:N	2.83	0.43
1:B:280:LEU:HD13	1:B:413:ILE:HG22	2.01	0.43
1:B:285:VAL:HG22	1:B:436:SER:HB3	2.01	0.43
1:B:439:LEU:HA	1:B:442:VAL:HG13	2.00	0.43
1:B:72:GLY:HA3	1:B:123:ASN:OD1	2.18	0.43
1:B:156:PHE:CD1	3:B:601:LMT:H72	2.54	0.43
1:B:371:PHE:CE1	1:B:470:LEU:HB3	2.54	0.43
1:A:401:GLY:O	1:A:405:VAL:HG22	2.18	0.42
1:A:487:LEU:HD23	1:A:487:LEU:HA	1.86	0.42
1:B:61:ILE:HG13	1:B:114:VAL:HG22	2.01	0.42
1:B:317:LEU:HD12	1:B:317:LEU:HA	1.92	0.42
1:A:226:LEU:HD23	1:A:226:LEU:HA	1.70	0.42
1:A:458:GLN:C	1:A:461:PRO:HD2	2.40	0.42
1:A:197:ALA:O	1:A:200:HIS:CB	2.67	0.42
1:B:115:VAL:HG22	1:B:230:SER:HB2	2.02	0.42
1:A:10:LYS:HB3	1:A:10:LYS:HE2	1.64	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:ARG:HA	1:A:241:GLU:OE2	2.20	0.42
1:A:82:MET:HE1	1:A:127:LEU:HD21	2.00	0.42
1:A:95:LEU:HA	1:A:95:LEU:HD12	1.64	0.42
1:A:16:THR:O	1:A:20:ILE:HG22	2.19	0.42
1:A:123:ASN:HA	1:A:126:ASN:HB2	2.01	0.42
1:B:96:MET:HG2	1:B:109:ALA:HB1	2.02	0.42
1:B:201:LYS:O	1:B:202:SER:C	2.57	0.42
1:A:374:TYR:CD1	1:A:390:VAL:HG12	2.55	0.42
1:A:472:VAL:HG22	1:B:372:PHE:CD1	2.54	0.42
1:B:34:LEU:HD22	1:B:165:LYS:HD3	2.01	0.42
1:A:378:GLY:O	1:A:380:PHE:N	2.53	0.42
1:A:407:GLY:O	1:A:412:MET:HG3	2.20	0.42
1:A:175:GLU:HA	1:A:176:PHE:O	2.20	0.41
1:A:317:LEU:HA	1:A:318:TRP:HA	1.74	0.41
1:A:156:PHE:CE1	3:A:602:LMT:H121	2.55	0.41
1:A:290:PHE:O	1:A:293:GLN:HB2	2.20	0.41
1:B:102:ASN:HD21	3:B:602:LMT:C3'	2.33	0.41
1:B:115:VAL:HG21	1:B:231:VAL:HG23	2.01	0.41
1:B:308:TRP:CE2	1:B:322:PRO:HD3	2.55	0.41
1:A:149:ALA:O	1:A:152:VAL:HG12	2.20	0.41
1:A:96:MET:SD	1:A:106:MET:HG3	2.60	0.41
1:A:376:PHE:O	1:A:376:PHE:CD1	2.74	0.41
1:A:439:LEU:HA	1:A:442:VAL:HG13	2.01	0.41
1:A:481:ALA:O	1:A:484:VAL:HG12	2.21	0.41
1:B:119:LEU:HD23	1:B:119:LEU:HA	1.78	0.41
1:B:126:ASN:O	1:B:130:LYS:HG2	2.21	0.41
1:A:37:TYR:HE2	1:A:323:ALA:N	2.18	0.41
3:A:602:LMT:H72	3:A:602:LMT:H102	1.61	0.41
1:B:197:ALA:O	1:B:200:HIS:CB	2.69	0.41
1:A:176:PHE:O	1:A:178:TRP:N	2.54	0.41
1:B:366:VAL:O	1:B:397:SER:OG	2.38	0.41
1:A:165:LYS:HA	1:A:177:GLY:O	2.20	0.41
1:A:256:LEU:O	1:A:260:PHE:CD2	2.74	0.41
1:A:356:ILE:HG12	1:A:356:ILE:O	2.21	0.41
1:B:95:LEU:HB3	1:B:109:ALA:HB2	2.02	0.41
1:B:140:ASP:O	1:B:144:THR:HG23	2.21	0.41
1:B:144:THR:O	1:B:147:TYR:HB3	2.21	0.41
1:B:163:TRP:HA	1:B:166:ASP:HB2	2.03	0.41
1:A:59:ALA:O	1:A:63:VAL:HG13	2.20	0.41
1:A:378:GLY:C	1:A:380:PHE:N	2.75	0.41
1:B:70:TRP:HE1	1:B:74:LYS:HD3	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:LEU:CD1	1:B:318:TRP:HA	2.48	0.41
3:A:602:LMT:O3'	3:A:602:LMT:H1B	2.21	0.40
1:B:49:ARG:HA	1:B:241:GLU:OE2	2.21	0.40
1:A:279:ILE:HG21	1:A:484:VAL:HG21	2.03	0.40
1:A:333:ILE:HG13	1:A:399:SER:HB2	2.04	0.40
1:A:485:LEU:N	1:A:486:PRO:HD2	2.36	0.40
1:B:378:GLY:HA2	1:B:382:VAL:HG21	2.02	0.40
1:A:410:LEU:H	1:A:410:LEU:HD12	1.87	0.40
1:B:332:TRP:CZ3	1:B:335:VAL:HG11	2.57	0.40
1:A:17:VAL:HG11	1:A:128:VAL:HG22	2.04	0.40
1:A:24:GLU:HA	1:A:189:ILE:HD13	2.03	0.40
1:A:39:MET:HE2	1:A:39:MET:HB3	1.84	0.40
1:A:293:GLN:HG3	1:A:470:LEU:CD1	2.51	0.40
1:A:311:GLN:NE2	1:A:312:VAL:C	2.75	0.40
1:A:8:VAL:HG11	1:A:12:HIS:CD2	2.57	0.40
3:A:602:LMT:H2'	3:A:602:LMT:H12	1.49	0.40
1:B:23:TRP:O	1:B:26:PHE:HB3	2.21	0.40
1:B:277:ALA:HB1	1:B:413:ILE:HD12	2.04	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	443/523 (85%)	403 (91%)	35 (8%)	5 (1%)	14	48
1	B	443/523 (85%)	403 (91%)	35 (8%)	5 (1%)	14	48
All	All	886/1046 (85%)	806 (91%)	70 (8%)	10 (1%)	14	48

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	205	ASN
1	A	208	SER
1	A	313	PHE
1	B	205	ASN
1	B	208	SER
1	A	387	SER
1	B	313	PHE
1	B	387	SER
1	B	209	GLU
1	A	209	GLU

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	340/417 (82%)	323 (95%)	17 (5%)	24	57
1	B	340/417 (82%)	331 (97%)	9 (3%)	46	74
All	All	680/834 (82%)	654 (96%)	26 (4%)	33	65

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LYS
1	A	43	LEU
1	A	90	SER
1	A	113	ILE
1	A	121	LYS
1	A	127	LEU
1	A	129	ARG
1	A	156	PHE
1	A	217	LYS
1	A	261	HIS
1	A	315	THR
1	A	337	SER
1	A	360	PHE
1	A	376	PHE

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Mol	Chain	Res	Type
1	A	380	PHE
1	A	489	ARG
1	A	490	ARG
1	B	13	SER
1	B	43	LEU
1	B	74	LYS
1	B	105	PHE
1	B	113	ILE
1	B	165	LYS
1	B	283	GLN
1	B	369	ILE
1	B	490	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	HIS
1	A	283	GLN
1	A	311	GLN
1	A	329	ASN
1	A	383	ASN
1	B	102	ASN
1	B	283	GLN
1	B	451	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 monosaccharides in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 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
5	DBY	A	604	4	14,15,15	1.08	1 (7%)	20,21,21	1.37	3 (15%)
3	LMT	B	602	-	36,36,36	1.23	5 (13%)	47,47,47	1.38	7 (14%)
4	ALA	A	603	5	3,4,5	1.09	0	2,4,6	1.24	0
3	LMT	A	602	-	36,36,36	1.27	3 (8%)	47,47,47	1.71	8 (17%)
3	LMT	B	601	-	36,36,36	1.17	2 (5%)	47,47,47	1.50	8 (17%)

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	DBY	A	604	4	-	3/8/8/8	0/1/1/1
3	LMT	B	602	-	-	11/21/61/61	0/2/2/2
4	ALA	A	603	5	-	0/0/2/4	-
3	LMT	A	602	-	-	12/21/61/61	0/2/2/2
3	LMT	B	601	-	-	14/21/61/61	0/2/2/2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	602	LMT	O5B-C1B	3.53	1.50	1.41
3	B	601	LMT	O5B-C1B	3.48	1.50	1.41
3	A	602	LMT	O5'-C1'	3.36	1.50	1.41
3	A	602	LMT	O5B-C1B	2.93	1.49	1.41
3	B	601	LMT	O5'-C1'	2.90	1.49	1.41
3	B	602	LMT	O5'-C1'	2.64	1.48	1.41
5	A	604	DBY	BR1-CE1	2.38	1.95	1.89
3	A	602	LMT	O5'-C5'	2.26	1.49	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	602	LMT	O5'-C5'	2.21	1.49	1.44
3	B	602	LMT	O1B-C4'	2.15	1.49	1.43
3	B	602	LMT	O3'-C3'	2.03	1.47	1.43

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	LMT	C1-O1'-C1'	4.99	122.11	113.84
3	A	602	LMT	C1'-C2'-C3'	4.41	119.18	110.00
3	B	601	LMT	C1'-C2'-C3'	4.40	119.15	110.00
3	A	602	LMT	C2'-C3'-C4'	4.05	118.94	109.68
3	B	602	LMT	C1-O1'-C1'	3.83	120.19	113.84
3	B	601	LMT	O5'-C5'-C6'	3.40	114.90	106.44
3	A	602	LMT	C1B-O1B-C4'	-3.24	109.95	117.96
3	B	602	LMT	O1B-C4'-C3'	3.16	115.69	107.28
3	A	602	LMT	O5'-C5'-C6'	3.03	113.97	106.44
3	B	601	LMT	O5B-C1B-C2B	2.84	116.36	110.35
3	B	602	LMT	O1'-C1'-C2'	2.83	112.72	108.30
3	B	601	LMT	C2'-C3'-C4'	2.80	116.07	109.68
3	B	601	LMT	C1-O1'-C1'	2.78	118.45	113.84
3	A	602	LMT	O5'-C1'-C2'	2.78	116.23	110.35
3	B	601	LMT	C1B-O1B-C4'	-2.73	111.20	117.96
5	A	604	DBY	CE2-CZ-CE1	2.72	119.94	116.56
3	B	602	LMT	O5B-C1B-C2B	2.66	115.97	110.35
3	B	601	LMT	O5'-C1'-C2'	2.59	115.84	110.35
3	A	602	LMT	O5B-C5B-C4B	2.40	114.05	109.69
3	B	602	LMT	O5'-C5'-C6'	2.34	112.26	106.44
5	A	604	DBY	CD1-CE1-CZ	-2.34	119.36	121.90
5	A	604	DBY	BR2-CE2-CZ	2.24	121.31	118.80
3	B	602	LMT	O4'-C4B-C3B	2.11	115.23	110.35
3	A	602	LMT	O3'-C3'-C2'	-2.10	105.49	110.35
3	B	602	LMT	O5'-C5'-C4'	2.03	114.03	109.75
3	B	601	LMT	C6'-C5'-C4'	-2.01	107.48	113.33

There are no chirality outliers.

All (40) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	602	LMT	C2'-C1'-O1'-C1
3	B	601	LMT	C2'-C1'-O1'-C1
3	B	601	LMT	O5'-C1'-O1'-C1
3	B	601	LMT	C2-C1-O1'-C1'

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Mol	Chain	Res	Type	Atoms
3	B	602	LMT	C2'-C1'-O1'-C1
3	B	601	LMT	O5'-C5'-C6'-O6'
3	A	602	LMT	C2-C3-C4-C5
3	B	602	LMT	O5'-C1'-O1'-C1
3	B	602	LMT	C4'-C5'-C6'-O6'
3	B	602	LMT	C3'-C4'-O1B-C1B
3	B	601	LMT	C4'-C5'-C6'-O6'
3	B	602	LMT	O5B-C5B-C6B-O6B
3	A	602	LMT	O1'-C1-C2-C3
3	B	602	LMT	C5'-C4'-O1B-C1B
3	B	601	LMT	C2-C3-C4-C5
3	A	602	LMT	C11-C10-C9-C8
3	B	602	LMT	C5-C6-C7-C8
3	A	602	LMT	C5-C6-C7-C8
3	B	601	LMT	C4-C5-C6-C7
3	A	602	LMT	C2-C1-O1'-C1'
3	B	602	LMT	O5'-C5'-C6'-O6'
3	B	602	LMT	C1-C2-C3-C4
3	A	602	LMT	C4-C5-C6-C7
3	B	601	LMT	O1'-C1-C2-C3
3	A	602	LMT	C3-C4-C5-C6
3	A	602	LMT	C7-C8-C9-C10
3	B	601	LMT	C9-C10-C11-C12
3	A	602	LMT	C3'-C4'-O1B-C1B
3	B	601	LMT	C1-C2-C3-C4
3	B	601	LMT	C5-C6-C7-C8
3	A	602	LMT	C5'-C4'-O1B-C1B
3	B	602	LMT	C3-C4-C5-C6
3	B	602	LMT	C4B-C5B-C6B-O6B
3	B	601	LMT	O5B-C5B-C6B-O6B
3	A	602	LMT	O5'-C1'-O1'-C1
5	A	604	DBY	O-C-CA-CB
5	A	604	DBY	OXT-C-CA-CB
3	B	601	LMT	C5'-C4'-O1B-C1B
3	B	601	LMT	C3'-C4'-O1B-C1B
5	A	604	DBY	OXT-C-CA-N

There are no ring outliers.

4 monomers are involved in 16 short contacts:

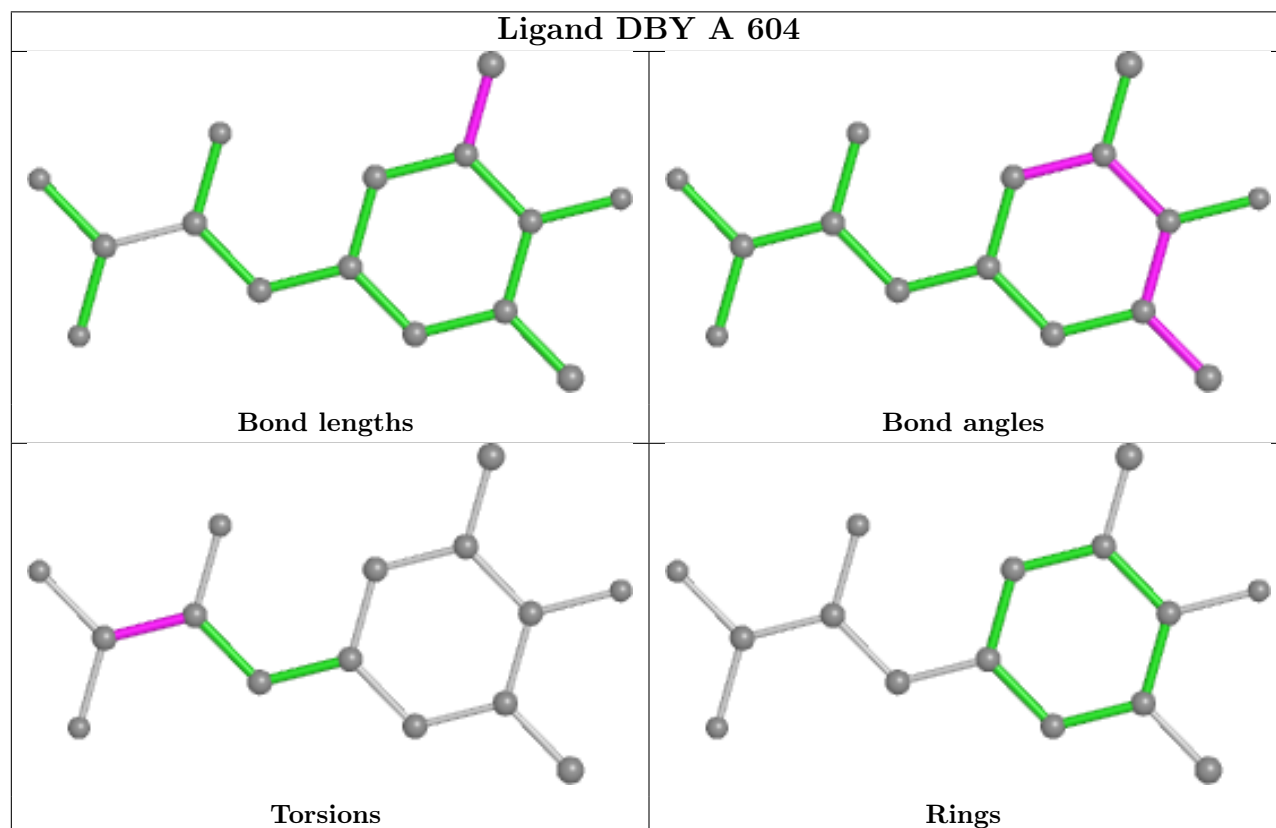
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	604	DBY	1	0

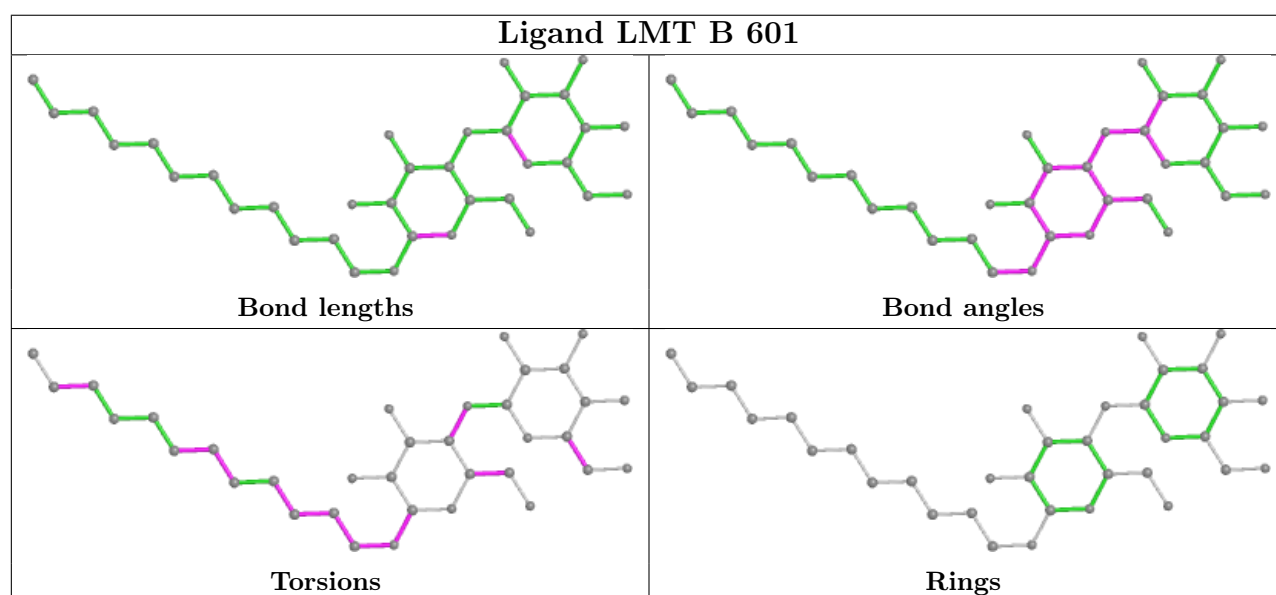
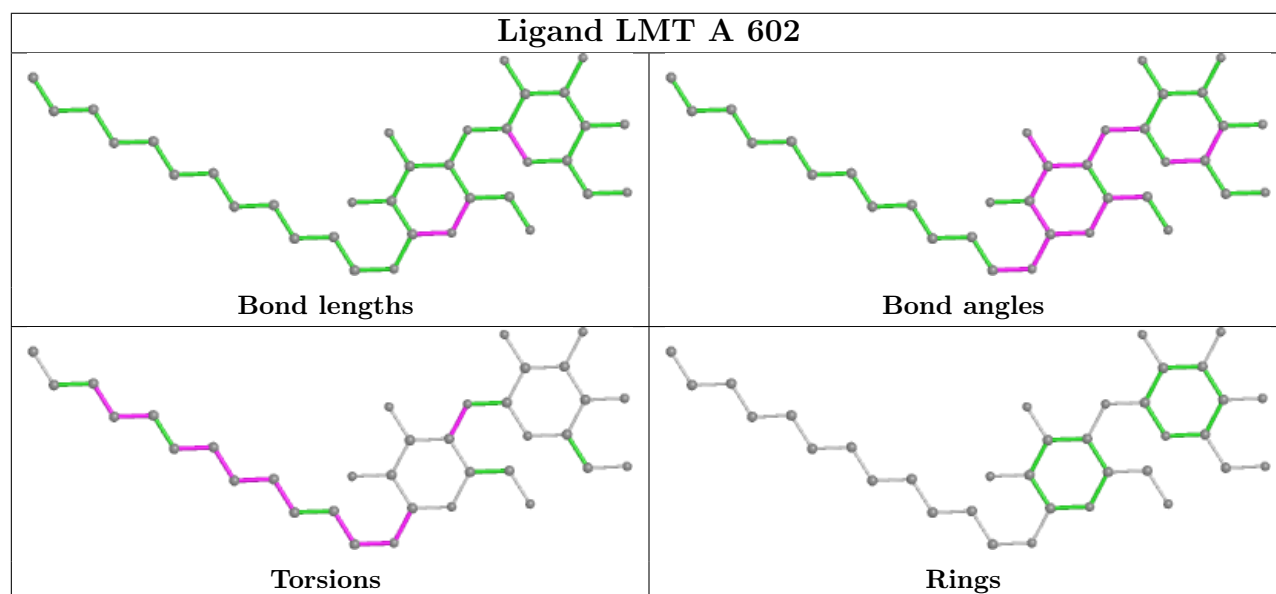
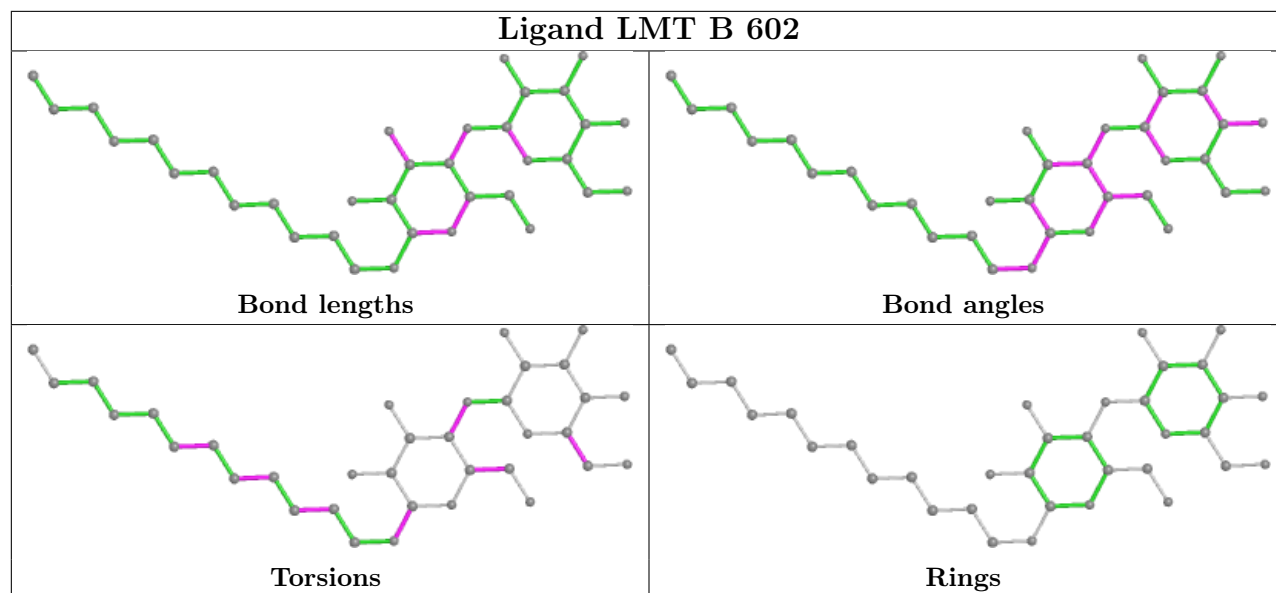
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	602	LMT	3	0
3	A	602	LMT	9	0
3	B	601	LMT	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	453/523 (86%)	-0.13	26 (5%) 23 12	61, 93, 141, 159	0
1	B	453/523 (86%)	-0.18	14 (3%) 49 32	56, 91, 143, 165	0
All	All	906/1046 (86%)	-0.16	40 (4%) 34 20	56, 93, 142, 165	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	315	THR	5.8
1	B	315	THR	5.7
1	A	312	VAL	5.1
1	B	202	SER	4.9
1	A	411	ALA	4.2
1	B	316	HIS	4.1
1	A	317	LEU	4.0
1	B	411	ALA	3.7
1	A	133	GLU	3.4
1	A	316	HIS	3.4
1	B	314	GLY	3.4
1	B	270	GLU	3.1
1	A	455	ASP	3.0
1	B	313	PHE	3.0
1	B	491	LEU	3.0
1	B	175	GLU	2.8
1	A	167	TYR	2.8
1	A	342	TRP	2.8
1	A	385	LYS	2.6
1	A	202	SER	2.6
1	A	270	GLU	2.5
1	A	386	THR	2.5
1	A	311	GLN	2.5
1	A	410	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	214	PRO	2.4
1	A	172	TYR	2.3
1	A	213	ARG	2.3
1	A	171	GLN	2.3
1	A	458	GLN	2.3
1	B	74	LYS	2.3
1	B	492	THR	2.3
1	B	317	LEU	2.2
1	B	312	VAL	2.2
1	A	313	PHE	2.2
1	A	426	MET	2.2
1	B	272	ALA	2.1
1	A	197	ALA	2.1
1	A	490	ARG	2.1
1	A	272	ALA	2.0
1	A	454	VAL	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](#)

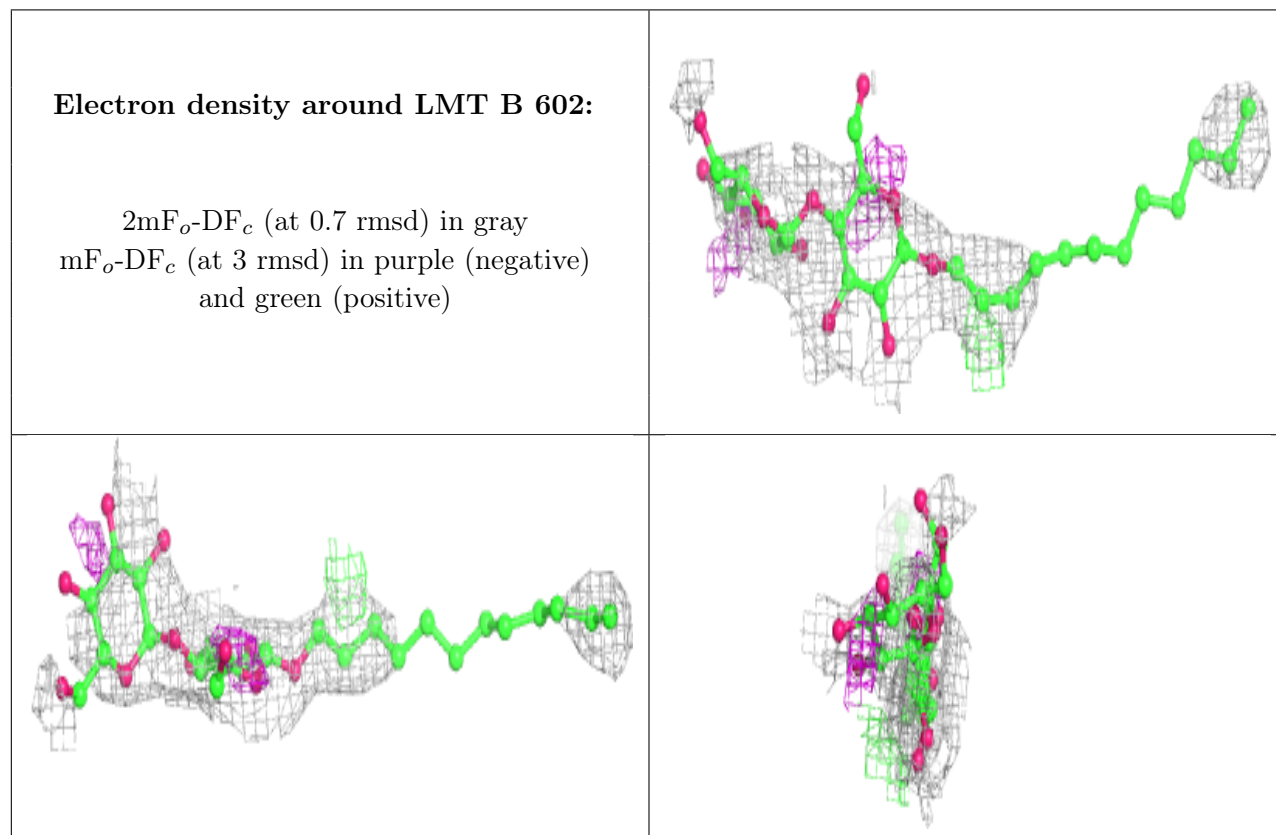
There are no monosaccharides in this entry.

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.

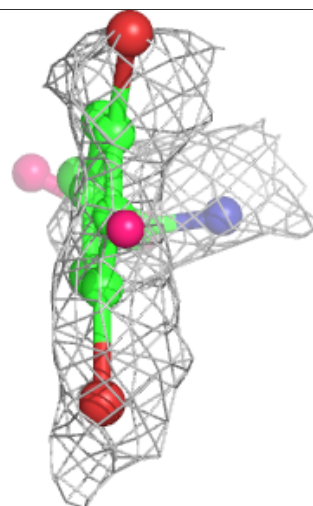
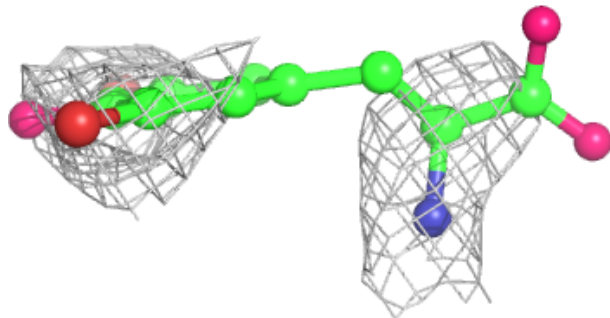
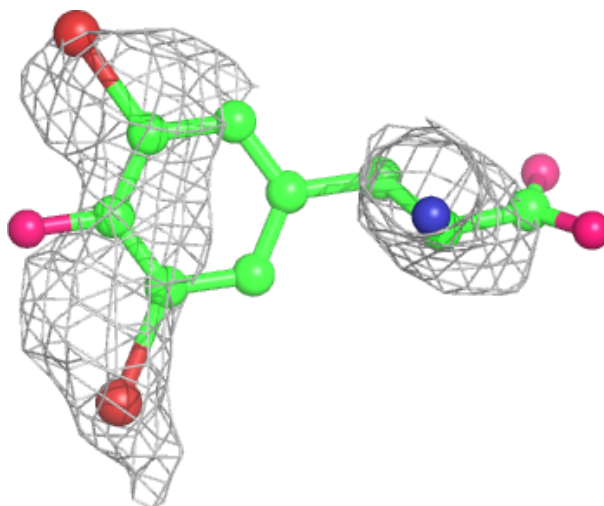
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	LMT	B	602	35/35	0.70	0.62	64,117,125,128	0
4	ALA	A	603	5/6	0.72	0.42	115,116,136,137	0
2	ZN	A	601	1/1	0.73	0.34	150,150,150,150	0
5	DBY	A	604	15/15	0.78	0.36	140,168,246,276	0
3	LMT	A	602	35/35	0.83	0.43	57,114,125,135	0
3	LMT	B	601	35/35	0.84	0.40	85,115,126,134	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.



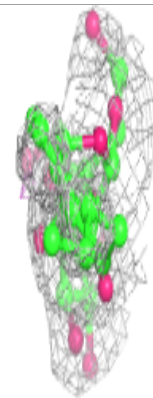
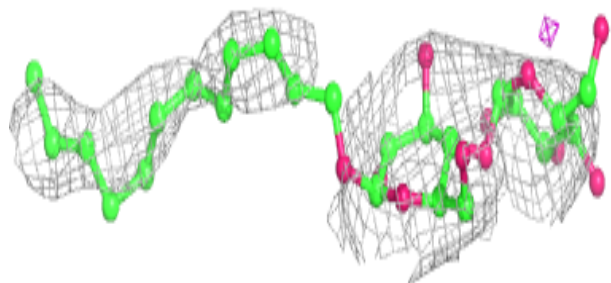
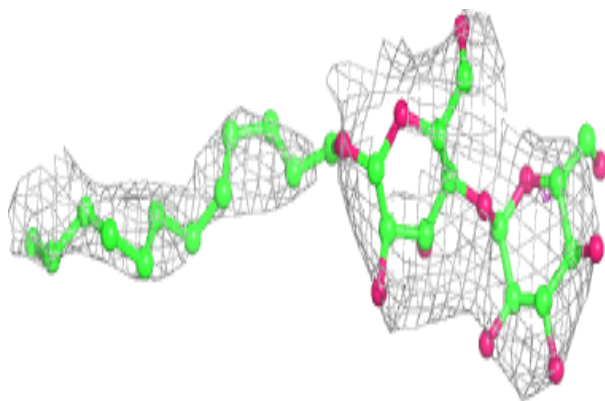
Electron density around DBY A 604:

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

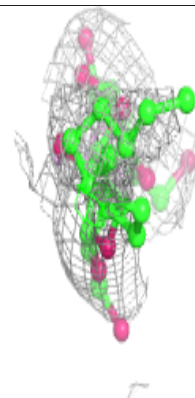
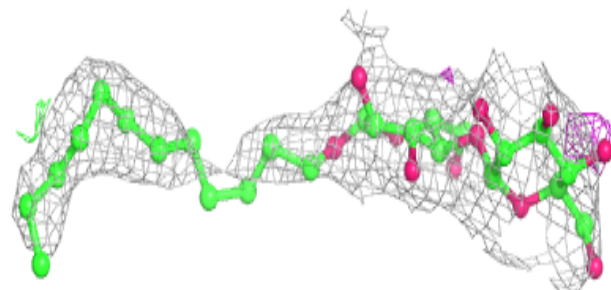
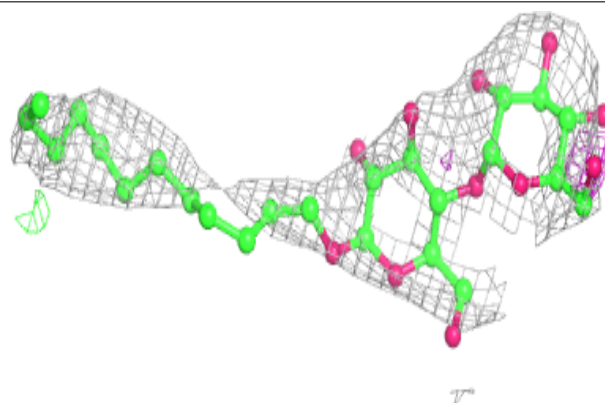


Electron density around LMT A 602:

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

**Electron density around LMT B 601:**

$2mF_o-DF_c$ (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.