



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:07 PM UTC

PDB ID : 7JS7 / pdb_00007js7
Title : Phospholipase D engineered mutant (TNYR) H442 covalent adduct with 1-
inositol phosphate
Authors : Vrieling, A.; Samantha, A.
Deposited on : 2020-08-13
Resolution : 2.30 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)
Xtrriage (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

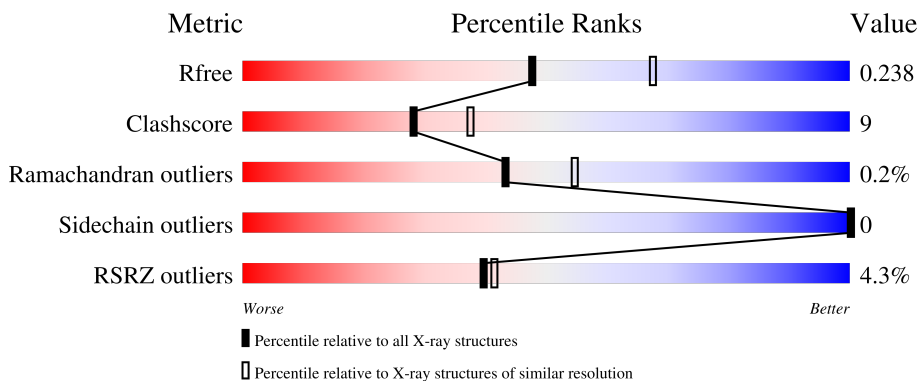
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	509	 4% 83% 14% ..
1	B	509	 5% 79% 17% .

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7806 atoms, of which 76 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 Phospholipase D.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	494	3705	2295	6	658	733	13	0	0	0
1	B	491	3679	2280	6	651	729	13	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

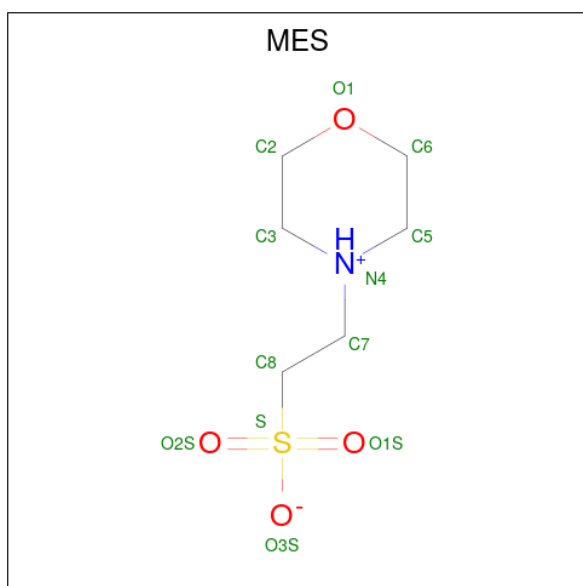
Chain	Residue	Modelled	Actual	Comment	Reference
A	107	SER	VAL	engineered mutation	UNP Q53728
A	186	THR	GLY	engineered mutation	UNP Q53728
A	187	ASN	TRP	engineered mutation	UNP Q53728
A	211	ARG	ALA	engineered mutation	UNP Q53728
A	377	ALA	ARG	conflict	UNP Q53728
A	385	ARG	TYR	engineered mutation	UNP Q53728
B	107	SER	VAL	engineered mutation	UNP Q53728
B	186	THR	GLY	engineered mutation	UNP Q53728
B	187	ASN	TRP	engineered mutation	UNP Q53728
B	211	ARG	ALA	engineered mutation	UNP Q53728
B	377	ALA	ARG	conflict	UNP Q53728
B	385	ARG	TYR	engineered mutation	UNP Q53728

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			14	3	8	3		
2	B	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 3 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



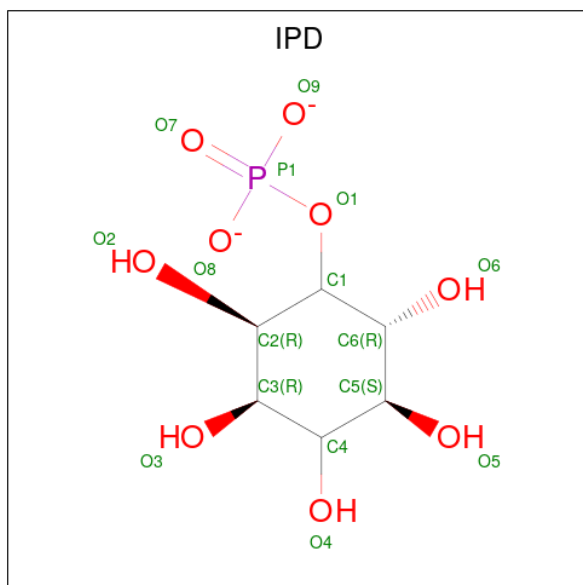
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
3	A	1	Total	C	H	N	O	S	0	0
			25	6	13	1	4	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	H	N	O			S
3	B	1	25	6	13	1	4	1	0	0

- Molecule 4 is D-MYO-INOSITOL-1-PHOSPHATE (CCD ID: IPD) (formula: C₆H₁₁O₉P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	H	O	P		
4	A	1	26	6	11	8	1	0	0
4	B	1	26	6	11	8	1	0	0

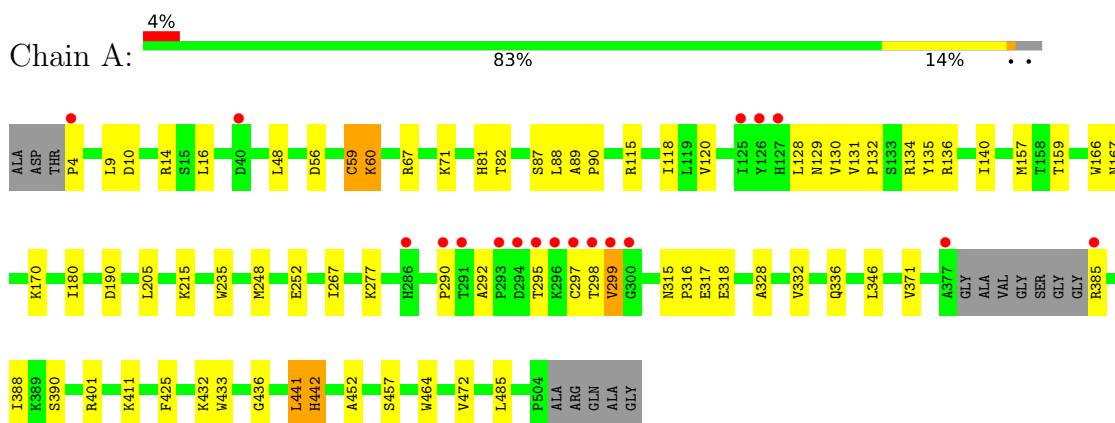
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	155	155	155	0	0
5	B	137	137	137	0	0

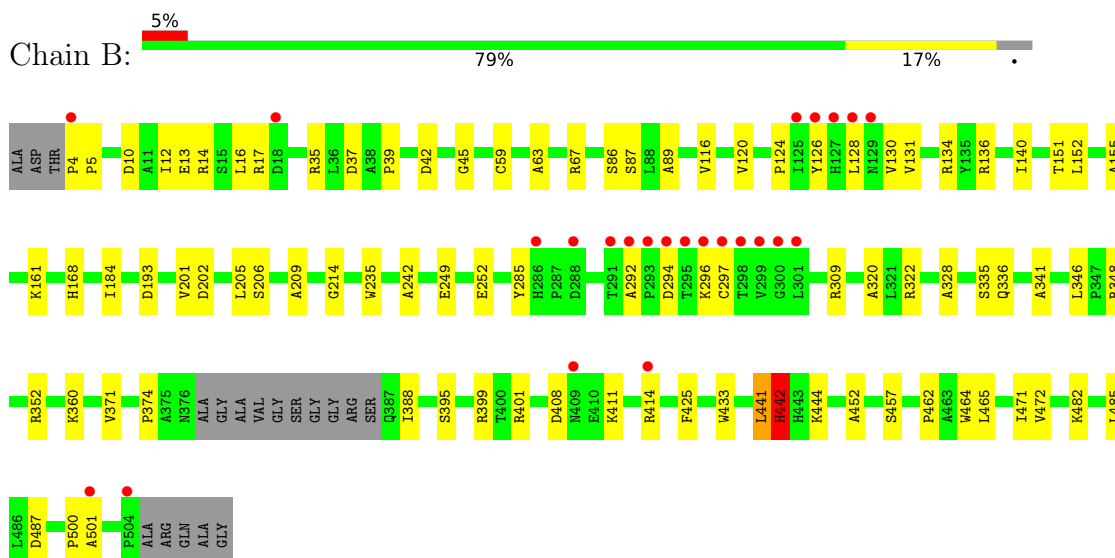
3 Residue-property plots [i](#)

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: Phospholipase D



- Molecule 1: Phospholipase D



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	90.58Å 98.81Å 105.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.29 – 2.30 45.29 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (45.29-2.30) 99.8 (45.29-2.30)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.83 (at 2.29Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660, PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.182 , 0.238 0.184 , 0.238	Depositor DCC
R_{free} test set	2100 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	16.8	Xtrriage
Anisotropy	0.919	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 36.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7806	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 48.81 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.1540e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: MES, GOL, IPD

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	0.40	2/3780 (0.1%)	0.70	6/5158 (0.1%)
1	B	0.39	3/3754 (0.1%)	0.63	4/5125 (0.1%)
All	All	0.39	5/7534 (0.1%)	0.67	10/10283 (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 (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	442	HIS	C-N	7.48	1.43	1.33
1	B	442	HIS	C-N	7.45	1.43	1.33
1	A	441	LEU	C-N	6.26	1.43	1.33
1	B	441	LEU	C-N	5.20	1.43	1.34
1	B	442	HIS	CG-ND1	-5.02	1.32	1.38

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	441	LEU	O-C-N	-20.90	98.92	123.16
1	A	441	LEU	O-C-N	-19.10	101.68	123.27
1	A	441	LEU	CA-C-N	15.14	146.16	122.08
1	A	441	LEU	C-N-CA	15.14	146.16	122.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	442	HIS	O-C-N	-9.71	109.96	122.20
1	B	441	LEU	CA-C-N	8.44	136.77	121.15
1	B	441	LEU	C-N-CA	8.44	136.77	121.15
1	B	442	HIS	O-C-N	-6.91	112.76	122.15
1	A	59	CYS	CA-CB-SG	-5.66	101.38	114.40
1	A	60	LYS	CB-CG-CD	-5.38	98.92	111.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	442	HIS	Mainchain
1	B	442	HIS	Mainchain

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	3699	6	3605	65	3
1	B	3673	6	3571	66	3
2	A	6	8	8	0	0
2	B	6	8	8	0	0
3	A	12	13	13	1	0
3	B	12	13	13	2	0
4	A	15	11	11	3	0
4	B	15	11	11	0	0
5	A	155	0	0	17	0
5	B	137	0	0	13	0
All	All	7730	76	7240	131	3

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 9.

All (131) 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:201:VAL:HB	5:B:705:HOH:O	1.70	0.90
1:B:13:GLU:OE2	1:B:17:ARG:NH1	2.09	0.86
1:A:118:ILE:HG22	5:A:762:HOH:O	1.77	0.83
1:A:120:VAL:HG13	5:A:762:HOH:O	1.77	0.83
1:B:131:VAL:HA	5:B:827:HOH:O	1.80	0.82
1:B:126:TYR:HE1	1:B:128:LEU:HD23	1.45	0.81
1:A:298:THR:OG1	5:A:701:HOH:O	2.00	0.79
1:B:130:VAL:HG11	5:B:783:HOH:O	1.83	0.79
1:A:115:ARG:HH21	1:A:115:ARG:HG3	1.48	0.77
1:B:130:VAL:HG23	1:B:134:ARG:HD3	1.70	0.74
1:A:59:CYS:HB2	1:A:433:TRP:CZ3	2.23	0.72
1:A:315:ASN:HB3	1:A:318:GLU:OE1	1.88	0.72
1:A:128:LEU:CD2	1:A:131:VAL:HA	2.19	0.72
1:B:482:LYS:HE3	5:B:701:HOH:O	1.88	0.71
1:A:295:THR:HA	5:A:701:HOH:O	1.94	0.68
1:A:60:LYS:HD3	1:A:60:LYS:C	2.19	0.68
1:A:215:LYS:HE3	5:A:773:HOH:O	1.94	0.68
1:B:374:PRO:HG3	1:B:388:ILE:CD1	2.25	0.67
1:B:128:LEU:HB3	5:B:827:HOH:O	1.94	0.66
1:B:285:TYR:O	1:B:309:ARG:HD2	1.95	0.66
1:A:157:MET:HE2	1:A:167:ASN:H	1.61	0.66
1:A:317:GLU:OE2	5:A:703:HOH:O	2.15	0.65
1:B:12:ILE:HD13	1:B:471:ILE:HD13	1.77	0.65
1:B:35:ARG:NH1	1:B:37:ASP:HA	2.12	0.64
1:A:128:LEU:HD21	1:A:131:VAL:HA	1.78	0.63
1:B:374:PRO:HG3	1:B:388:ILE:HD11	1.80	0.63
1:A:88:LEU:CD2	4:A:603:IPD:H3	2.29	0.62
1:B:193:ASP:HB3	5:B:831:HOH:O	1.99	0.62
1:A:295:THR:HG22	1:A:295:THR:O	2.00	0.62
1:B:294:ASP:OD2	1:B:348:ARG:HD3	2.00	0.61
1:B:89:ALA:HB1	1:B:128:LEU:HD21	1.82	0.61
1:A:134:ARG:HD3	5:A:706:HOH:O	2.01	0.59
1:A:88:LEU:HD23	4:A:603:IPD:H3	1.84	0.59
1:B:10:ASP:O	1:B:14:ARG:HG3	2.04	0.57
1:B:126:TYR:CE1	1:B:128:LEU:HD23	2.33	0.56
1:B:87:SER:O	1:B:120:VAL:HA	2.05	0.56
1:A:67:ARG:HG3	5:A:755:HOH:O	2.03	0.56
1:B:130:VAL:CG1	5:B:754:HOH:O	2.53	0.56
1:B:336:GLN:O	1:B:371:VAL:HA	2.06	0.56
1:A:157:MET:HE1	1:A:159:THR:HG23	1.88	0.56
1:A:4:PRO:HD3	5:A:848:HOH:O	2.06	0.56
1:A:87:SER:O	1:A:120:VAL:HA	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:252:GLU:OE2	1:A:252:GLU:HA	2.06	0.55
1:B:16:LEU:HD21	1:B:320:ALA:HB2	1.87	0.55
1:B:205:LEU:HD23	1:B:206:SER:N	2.22	0.54
1:B:408:ASP:OD1	1:B:411:LYS:HE2	2.08	0.54
1:B:63:ALA:O	1:B:67:ARG:HG2	2.09	0.53
1:B:296:LYS:O	1:B:296:LYS:HE3	2.09	0.53
1:B:39:PRO:HD2	1:B:45:GLY:HA3	1.89	0.52
1:B:130:VAL:HG12	5:B:754:HOH:O	2.10	0.52
1:A:128:LEU:HD23	1:A:131:VAL:HA	1.90	0.52
1:A:425:PHE:HB2	1:A:441:LEU:HD22	1.91	0.52
1:B:346:LEU:HD13	1:B:464:TRP:CZ2	2.45	0.52
1:A:432:LYS:HD2	1:A:436:GLY:HA2	1.91	0.52
1:A:292:ALA:HB2	1:A:401:ARG:NH1	2.25	0.51
1:A:136:ARG:HD2	1:A:140:ILE:HD11	1.93	0.50
1:A:290:PRO:HB3	5:A:845:HOH:O	2.10	0.50
1:B:86:SER:HB2	5:B:766:HOH:O	2.12	0.49
1:B:485:LEU:C	1:B:485:LEU:HD23	2.37	0.49
1:B:252:GLU:HB2	5:B:702:HOH:O	2.11	0.49
1:A:48:LEU:HA	5:A:795:HOH:O	2.12	0.49
1:B:130:VAL:HG13	1:B:130:VAL:O	2.12	0.49
1:A:157:MET:CE	1:A:159:THR:HG23	2.42	0.49
1:A:157:MET:HE2	1:A:167:ASN:N	2.27	0.48
1:B:39:PRO:HG2	1:B:42:ASP:HB3	1.93	0.48
1:A:56:ASP:OD2	1:A:60:LYS:HB3	2.13	0.48
1:B:414:ARG:HG2	1:B:414:ARG:HH11	1.78	0.48
1:B:296:LYS:HE2	1:B:348:ARG:HH11	1.79	0.48
1:A:128:LEU:HD23	1:A:131:VAL:HG12	1.96	0.47
1:A:157:MET:CE	1:A:166:TRP:HA	2.44	0.47
1:A:315:ASN:O	1:A:318:GLU:OE1	2.33	0.47
1:B:128:LEU:HD23	1:B:128:LEU:HA	1.72	0.46
1:A:88:LEU:HD21	4:A:603:IPD:H3	1.97	0.46
1:A:170:LYS:HD2	1:A:457:SER:HB3	1.98	0.46
1:B:399:ARG:NH2	1:B:501:ALA:O	2.45	0.46
1:B:151:THR:HG23	1:B:242:ALA:HB2	1.96	0.46
1:B:444:LYS:HD3	1:B:457:SER:HA	1.98	0.45
1:A:9:LEU:HD22	1:A:267:ILE:HD12	1.97	0.45
1:B:360:LYS:HD2	1:B:360:LYS:HA	1.74	0.45
1:A:128:LEU:HD21	1:A:132:PRO:HD3	1.98	0.45
1:A:388:ILE:HD12	1:A:390:SER:O	2.16	0.45
1:B:395:SER:HB2	1:B:500:PRO:HB3	1.98	0.45
1:B:59:CYS:HB2	1:B:433:TRP:CZ3	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:452:ALA:HA	1:A:472:VAL:O	2.17	0.45
1:B:205:LEU:HD21	1:B:209:ALA:HB3	1.98	0.45
1:B:335:SER:C	1:B:336:GLN:HG2	2.42	0.45
1:A:205:LEU:C	1:A:205:LEU:HD12	2.42	0.45
1:A:298:THR:O	1:A:299:VAL:C	2.60	0.45
1:A:277:LYS:NZ	5:A:702:HOH:O	2.10	0.44
1:A:130:VAL:HG12	1:A:135:TYR:HB2	1.98	0.44
1:A:4:PRO:HD3	5:A:815:HOH:O	2.17	0.44
1:B:487:ASP:OD2	5:B:701:HOH:O	2.21	0.44
1:A:485:LEU:C	1:A:485:LEU:HD23	2.43	0.44
1:B:184:ILE:HD12	1:B:202:ASP:CG	2.42	0.44
1:A:411:LYS:HD2	5:A:736:HOH:O	2.17	0.44
1:B:136:ARG:O	1:B:140:ILE:HG13	2.18	0.44
1:B:168:HIS:HB2	1:B:465:LEU:HD13	1.99	0.43
1:A:16:LEU:HD22	1:A:316:PRO:HB2	2.00	0.43
1:B:328:ALA:O	1:B:360:LYS:HE3	2.18	0.43
1:B:322:ARG:HD3	1:B:352:ARG:HB3	2.01	0.43
1:A:277:LYS:HD3	5:A:702:HOH:O	2.18	0.43
1:B:124:PRO:HB3	1:B:161:LYS:HD2	2.00	0.43
1:B:425:PHE:HB2	1:B:441:LEU:HD22	2.01	0.43
1:A:190:ASP:OD2	1:A:385:ARG:NH2	2.52	0.43
1:A:328:ALA:CB	1:A:332:VAL:HG22	2.48	0.42
1:B:4:PRO:HA	1:B:5:PRO:HD3	1.81	0.42
1:A:115:ARG:HH21	1:A:115:ARG:CG	2.24	0.42
1:A:129:ASN:HB2	5:A:829:HOH:O	2.18	0.42
1:B:452:ALA:HA	1:B:472:VAL:O	2.19	0.42
1:A:235:TRP:CE3	3:A:602:MES:H21	2.55	0.42
1:A:336:GLN:O	1:A:371:VAL:HA	2.20	0.42
1:B:205:LEU:HD23	1:B:205:LEU:C	2.44	0.42
1:B:296:LYS:O	1:B:297:CYS:HB2	2.19	0.42
1:B:235:TRP:CE2	3:B:602:MES:H52	2.55	0.41
1:B:442:HIS:HB2	5:B:759:HOH:O	2.20	0.41
1:A:346:LEU:HD13	1:A:464:TRP:CZ2	2.55	0.41
1:A:60:LYS:C	1:A:60:LYS:CD	2.90	0.41
1:A:290:PRO:HG3	5:A:845:HOH:O	2.21	0.41
1:B:131:VAL:HG12	5:B:827:HOH:O	2.21	0.41
1:B:214:GLY:O	1:B:249:GLU:HG3	2.21	0.41
1:A:10:ASP:O	1:A:14:ARG:HG3	2.21	0.41
1:A:81:HIS:HB3	1:A:115:ARG:HH12	1.85	0.41
1:A:89:ALA:HB1	1:A:90:PRO:HA	2.03	0.41
1:B:116:VAL:HB	1:B:152:LEU:HD23	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:LYS:HG3	1:A:180:ILE:HD11	2.03	0.41
1:B:235:TRP:CD2	3:B:602:MES:H52	2.55	0.40
1:B:292:ALA:HB2	1:B:401:ARG:NH1	2.36	0.40
1:B:341:ALA:HB2	1:B:462:PRO:HD2	2.03	0.40
1:B:155:ALA:HA	1:B:235:TRP:O	2.21	0.40
1:A:82:THR:HB	1:A:248:MET:HE3	2.04	0.40
1:A:295:THR:O	1:A:295:THR:CG2	2.69	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:LYS:NZ	1:B:17:ARG:NE[1_455]	1.34	0.86
1:A:60:LYS:NZ	1:B:17:ARG:CD[1_455]	1.59	0.61
1:A:60:LYS:NZ	1:B:17:ARG:CZ[1_455]	1.88	0.32

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	490/509 (96%)	469 (96%)	19 (4%)	2 (0%)	30	38
1	B	487/509 (96%)	468 (96%)	19 (4%)	0	100	100
All	All	977/1018 (96%)	937 (96%)	38 (4%)	2 (0%)	43	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	297	CYS
1	A	299	VAL

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	395/401 (98%)	395 (100%)	0	100	100
1	B	392/401 (98%)	392 (100%)	0	100	100
All	All	787/802 (98%)	787 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	GLN
1	A	129	ASN
1	A	337	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](#)

6 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
3	MES	B	602	-	12,12,12	1.73	3 (25%)	15,16,16	1.65	4 (26%)
4	IPD	B	603	1	13,15,16	1.08	1 (7%)	18,22,25	1.24	2 (11%)
2	GOL	B	601	-	5,5,5	0.57	0	5,5,5	0.37	0
4	IPD	A	603	1	13,15,16	1.42	1 (7%)	18,22,25	1.90	6 (33%)
2	GOL	A	601	-	5,5,5	0.58	0	5,5,5	0.22	0
3	MES	A	602	-	12,12,12	1.66	3 (25%)	15,16,16	1.79	4 (26%)

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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MES	B	602	-	-	0/6/14/14	0/1/1/1
4	IPD	B	603	1	-	0/2/28/29	0/1/1/1
2	GOL	B	601	-	-	2/4/4/4	-
4	IPD	A	603	1	-	0/2/28/29	0/1/1/1
2	GOL	A	601	-	-	2/4/4/4	-
3	MES	A	602	-	-	3/6/14/14	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	602	MES	C8-S	4.27	1.83	1.77
3	A	602	MES	C8-S	3.91	1.83	1.77
4	A	603	IPD	O1-C1	-3.81	1.38	1.45
4	B	603	IPD	O1-C1	-2.97	1.40	1.45
3	A	602	MES	O2S-S	2.21	1.51	1.45
3	A	602	MES	O1S-S	2.21	1.51	1.45
3	B	602	MES	O1S-S	2.14	1.51	1.45
3	B	602	MES	O2S-S	2.13	1.51	1.45

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	603	IPD	O1-C1-C2	4.61	114.51	107.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	602	MES	O3S-S-O2S	-3.74	102.05	111.40
3	A	602	MES	O3S-S-O2S	-3.59	102.41	111.40
3	A	602	MES	O1S-S-C8	3.05	111.33	106.73
4	A	603	IPD	C5-C4-C3	2.92	115.96	110.83
4	A	603	IPD	C3-C2-C1	-2.91	103.07	109.68
3	A	602	MES	C2-C3-N4	2.89	114.52	110.12
3	B	602	MES	O2S-S-C8	2.52	110.54	106.73
4	A	603	IPD	O3-C3-C4	-2.40	104.72	110.38
3	A	602	MES	C8-C7-N4	-2.30	103.66	112.36
3	B	602	MES	O1S-S-C8	2.28	110.17	106.73
4	B	603	IPD	C6-C5-C4	2.25	114.79	110.83
4	B	603	IPD	O1-C1-C2	2.17	111.01	107.90
4	A	603	IPD	O1-C1-C6	-2.13	104.84	107.90
3	B	602	MES	O3S-S-C8	2.12	110.16	106.00
4	A	603	IPD	O3-C3-C2	2.04	115.18	110.38

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	601	GOL	C1-C2-C3-O3
3	A	602	MES	C7-C8-S-O2S
2	B	601	GOL	O2-C2-C3-O3
3	A	602	MES	C7-C8-S-O3S
2	A	601	GOL	C1-C2-C3-O3
2	A	601	GOL	O2-C2-C3-O3
3	A	602	MES	C7-C8-S-O1S

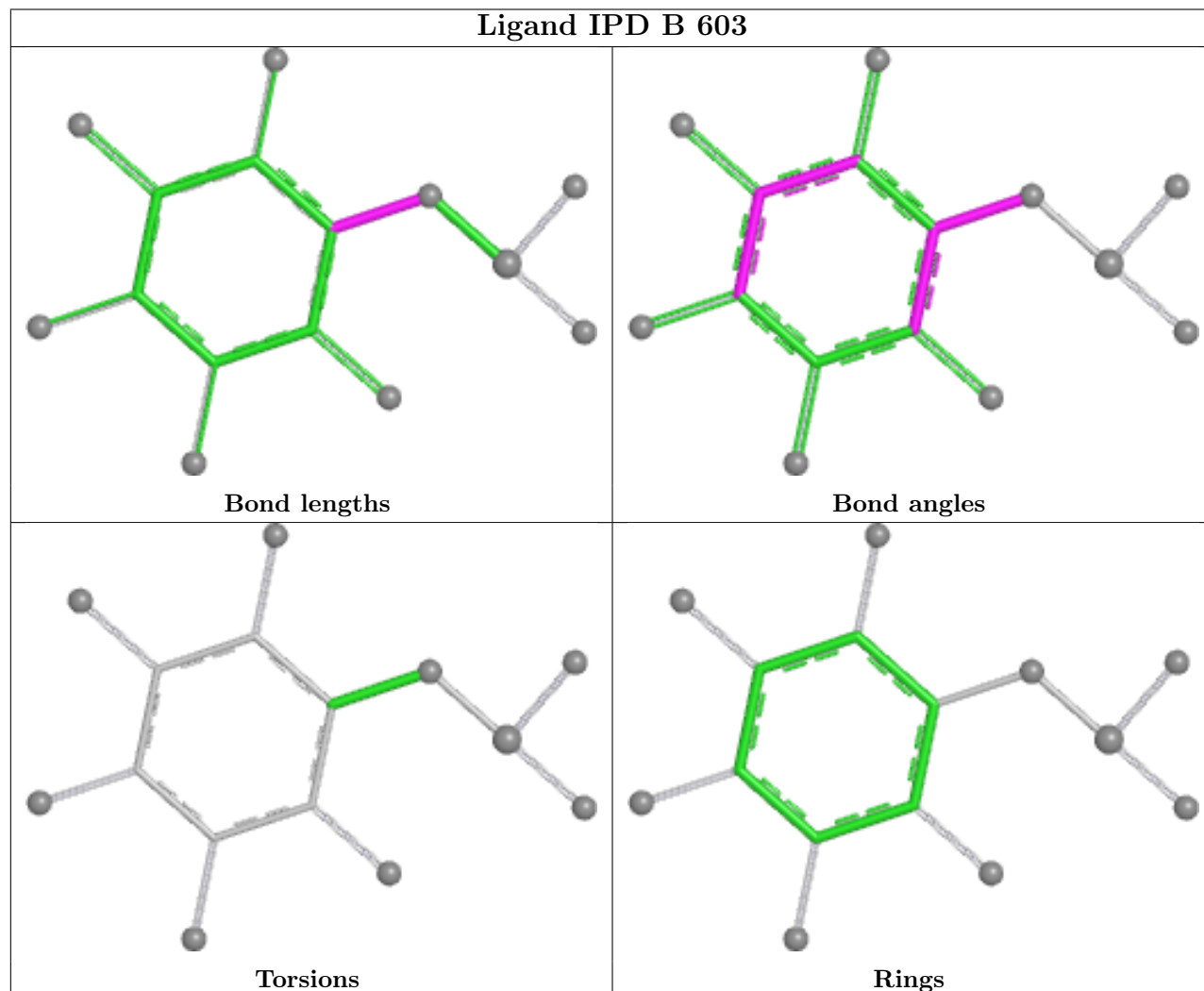
There are no ring outliers.

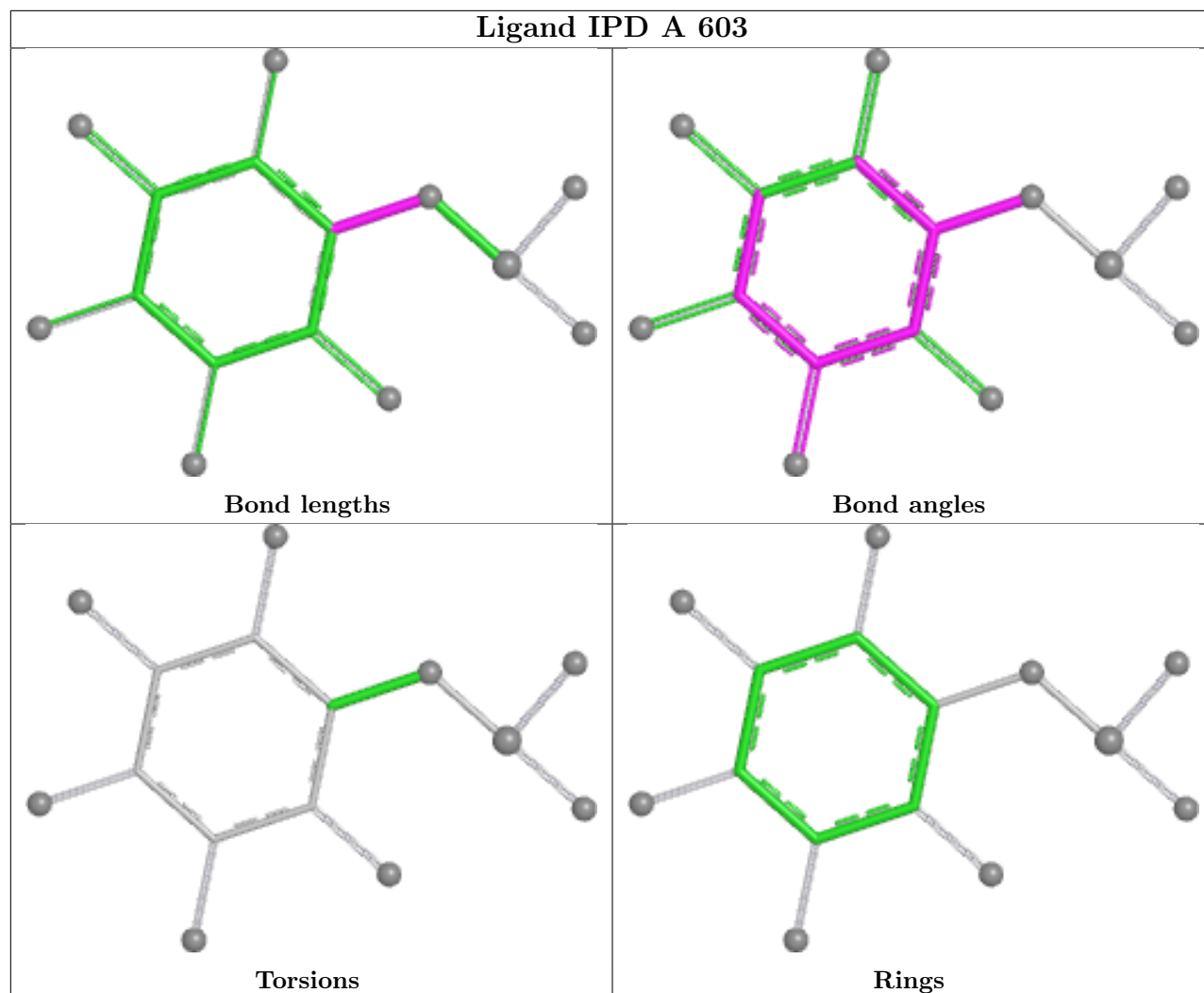
3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	602	MES	2	0
4	A	603	IPD	3	0
3	A	602	MES	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	A	494/509 (97%)	-0.16	18 (3%) 46 48	9, 17, 39, 90	4 (0%)
1	B	491/509 (96%)	-0.07	24 (4%) 35 36	10, 19, 47, 90	4 (0%)
All	All	985/1018 (96%)	-0.11	42 (4%) 40 41	9, 18, 41, 90	8 (0%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	295	THR	6.1
1	A	299	VAL	5.7
1	A	295	THR	5.7
1	A	298	THR	5.0
1	A	297	CYS	4.9
1	B	296	LYS	4.6
1	B	293	PRO	4.5
1	A	125	ILE	4.5
1	B	299	VAL	4.4
1	B	298	THR	4.2
1	B	300	GLY	3.9
1	B	501	ALA	3.9
1	B	127	HIS	3.9
1	A	296	LYS	3.7
1	B	291	THR	3.7
1	A	127	HIS	3.6
1	A	291	THR	3.4
1	B	292	ALA	3.3
1	A	294	ASP	3.2
1	B	297	CYS	3.2
1	A	300	GLY	3.2
1	A	293	PRO	3.1
1	A	4	PRO	3.0
1	B	286	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	18	ASP	2.7
1	B	288	ASP	2.6
1	B	128	LEU	2.5
1	B	504	PRO	2.5
1	A	290	PRO	2.5
1	B	409	ASN	2.5
1	B	126	TYR	2.4
1	A	385	ARG	2.4
1	A	286	HIS	2.4
1	B	294	ASP	2.3
1	B	301	LEU	2.2
1	A	126	TYR	2.2
1	B	125	ILE	2.2
1	B	4	PRO	2.1
1	B	414	ARG	2.1
1	B	129	ASN	2.1
1	A	377	ALA	2.0
1	A	40	ASP	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.

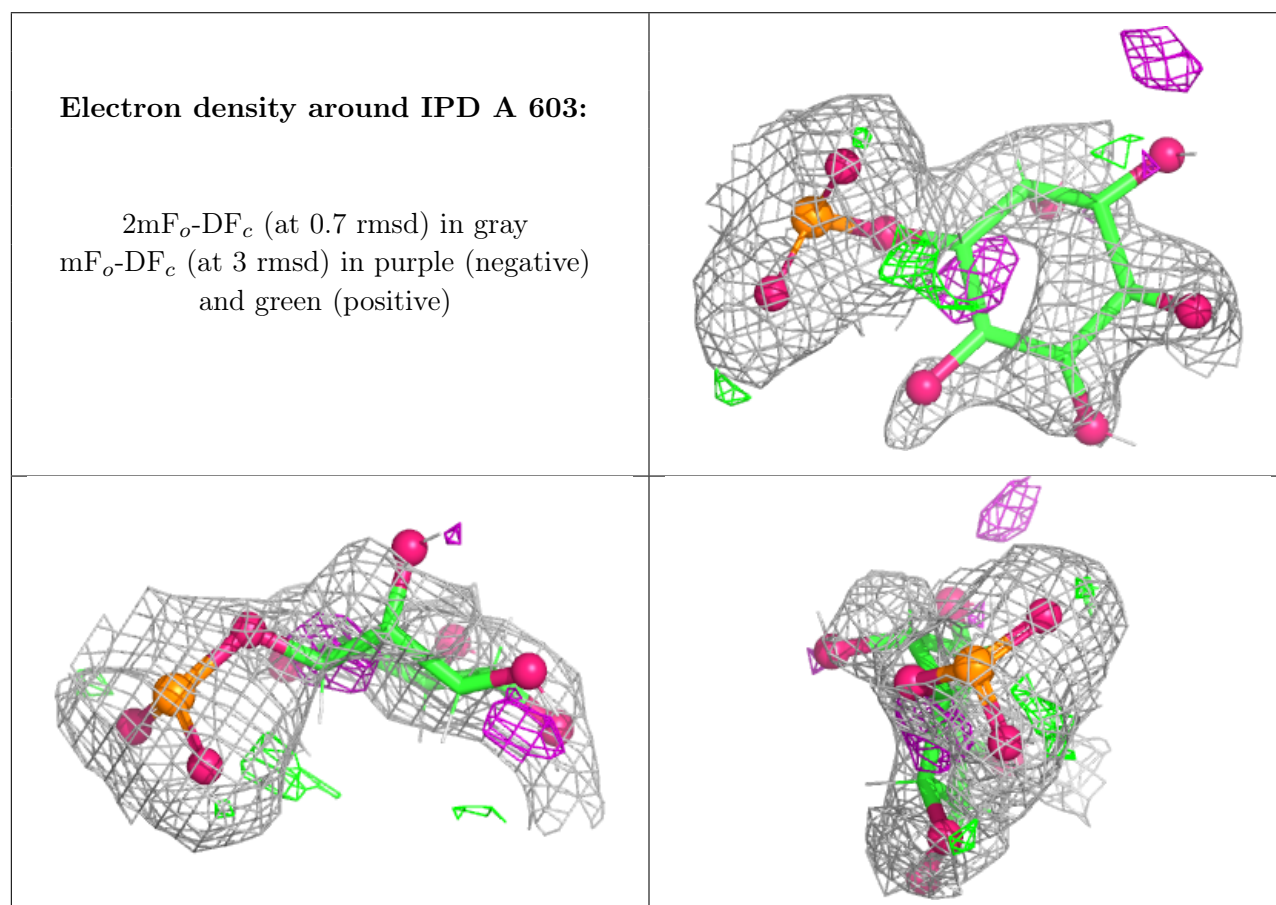
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	IPD	A	603	15/16	0.75	0.17	22,51,61,68	26
4	IPD	B	603	15/16	0.83	0.14	22,47,58,63	26
3	MES	A	602	12/12	0.87	0.11	23,33,50,51	0
3	MES	B	602	12/12	0.91	0.09	23,32,41,41	0
2	GOL	A	601	6/6	0.96	0.09	16,20,23,24	0

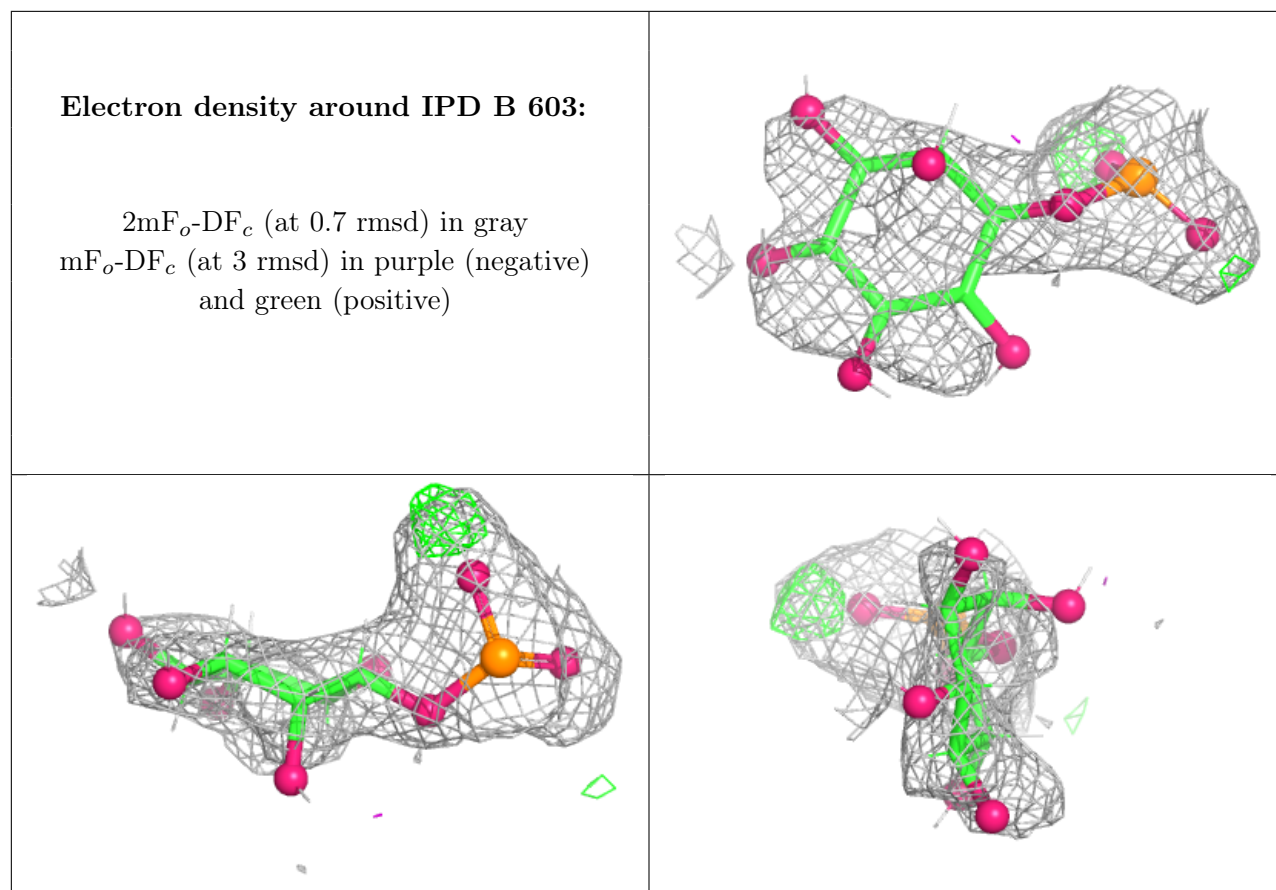
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	GOL	B	601	6/6	0.97	0.06	16,20,25,25	0

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.





6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.