



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 10:02 PM UTC

PDB ID : 2ZEU / pdb_00002zeu
Title : S. Cerevisiae Geranylgeranyl Pyrophosphate Synthase in Complex with BPH-715
Authors : Guo, R.T.; Chen, C.K.-M.; Cao, R.; Oldfield, E.; Wang, A.H.-J.
Deposited on : 2007-12-17
Resolution : 2.00 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

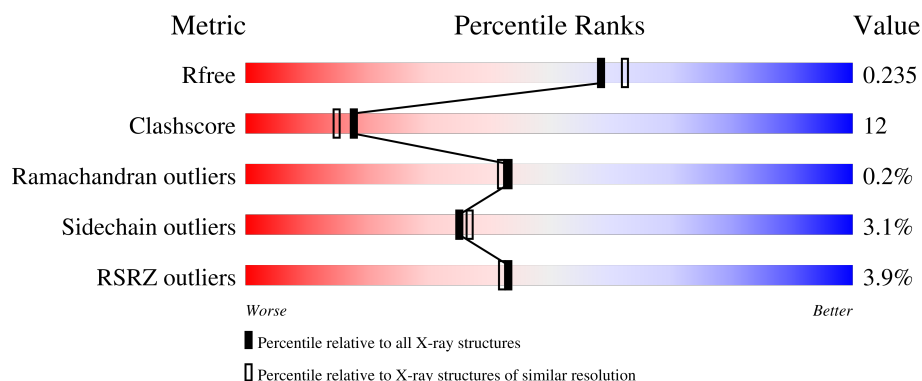
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	340	2% 69% 18% • 10%
1	B	340	5% 68% 19% • 9%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5614 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.

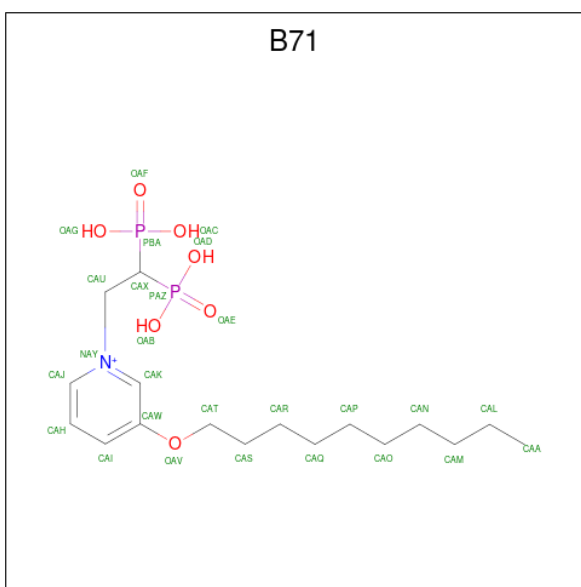
- Molecule 1 is a protein called Geranylgeranyl pyrophosphate synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	305	Total	C	N	O	S	0	0	0
			2491	1602	419	460	10			
1	B	308	Total	C	N	O	S	0	0	0
			2526	1625	424	468	9			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q12051
A	2	THR	-	expression tag	UNP Q12051
A	3	LYS	-	expression tag	UNP Q12051
A	4	ASN	-	expression tag	UNP Q12051
A	5	LYS	-	expression tag	UNP Q12051
B	1	MET	-	expression tag	UNP Q12051
B	2	THR	-	expression tag	UNP Q12051
B	3	LYS	-	expression tag	UNP Q12051
B	4	ASN	-	expression tag	UNP Q12051
B	5	LYS	-	expression tag	UNP Q12051

- Molecule 2 is 3-(DECYLOXY)-1-(2,2-DIPHOSPHONOETHYL)PYRIDINIUM (CCD ID: B71) (formula: C₁₇H₃₂NO₇P₂).

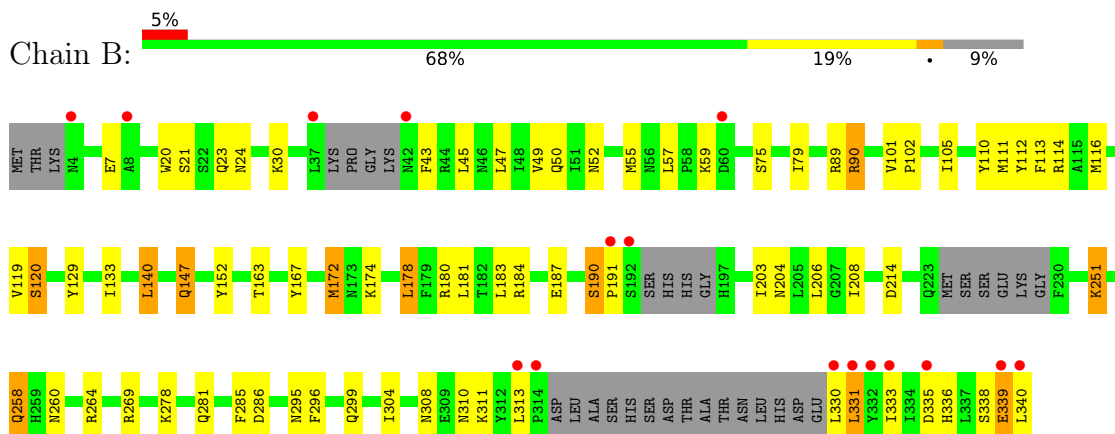


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 27	C 17	N 1	O 7	P 2	0	0
2	B	1	Total 27	C 17	N 1	O 7	P 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	283	Total O 283 283	0	0
3	B	260	Total O 260 260	0	0

- Molecule 1: Geranylgeranyl pyrophosphate synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	47.26Å 116.21Å 129.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.00 30.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.9 (30.00-2.00) 96.0 (30.00-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.54 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.189 , 0.236 0.189 , 0.235	Depositor DCC
R_{free} test set	2349 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	32.6	Xtriage
Anisotropy	0.741	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 49.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5614	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.58% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: B71

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	1.06	5/2533 (0.2%)	1.09	10/3424 (0.3%)
1	B	0.95	2/2570 (0.1%)	1.07	11/3478 (0.3%)
All	All	1.01	7/5103 (0.1%)	1.08	21/6902 (0.3%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	172	MET	SD-CE	-8.83	1.57	1.79
1	A	101	VAL	CA-CB	8.39	1.59	1.53
1	A	166	MET	SD-CE	-7.88	1.59	1.79
1	A	186	MET	SD-CE	6.45	1.95	1.79
1	A	170	MET	SD-CE	5.85	1.94	1.79
1	A	209	ILE	CA-CB	5.39	1.60	1.54
1	B	23	GLN	CA-C	5.01	1.59	1.52

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	59	LYS	N-CA-C	7.86	119.84	111.28
1	A	282	ILE	N-CA-C	-7.31	103.34	110.72
1	B	174	LYS	N-CA-C	7.07	118.68	110.97
1	B	331	LEU	N-CA-C	-6.76	103.00	111.11
1	B	191	PRO	N-CA-C	-6.10	108.33	114.68
1	A	174	LYS	N-CA-C	6.05	117.57	110.97
1	A	141	ILE	N-CA-C	-6.00	104.66	110.42
1	B	140	LEU	N-CA-C	-5.86	104.97	111.36
1	A	44	ARG	N-CA-C	-5.69	105.15	111.36
1	A	271	SER	N-CA-C	-5.66	105.58	112.88
1	A	105	ILE	CB-CA-C	-5.58	104.83	111.97
1	B	163	THR	N-CA-C	-5.42	102.44	110.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	140	LEU	N-CA-C	-5.40	105.42	112.23
1	A	238	LYS	N-CA-C	5.39	118.04	110.23
1	B	105	ILE	CB-CA-C	-5.31	105.17	111.97
1	A	59	LYS	N-CA-C	5.27	117.71	111.33
1	B	190	SER	N-CA-C	5.23	117.14	109.71
1	A	256	THR	N-CA-C	5.14	116.57	111.07
1	B	114	ARG	N-CA-C	-5.11	105.79	111.36
1	B	285	PHE	N-CA-C	5.06	119.53	112.90
1	B	269	ARG	N-CA-C	-5.05	104.42	111.39

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2491	0	2542	66	0
1	B	2526	0	2570	64	0
2	A	27	0	28	2	0
2	B	27	0	28	5	0
3	A	283	0	0	12	0
3	B	260	0	0	11	1
All	All	5614	0	5168	126	1

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 12.

All (126) 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:A:125:LYS:HE3	1:A:128:LEU:HB2	1.38	1.06
1:B:55:MET:HE2	1:B:187:GLU:HA	1.35	1.05
1:A:80:ASP:HB2	3:A:1401:HOH:O	1.62	0.99
1:A:52:ASN:HD21	1:A:57:LEU:H	1.10	0.97
1:B:172:MET:HE3	1:B:208:ILE:HD11	1.46	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:GLN:HE21	1:A:258:GLN:H	1.12	0.96
1:A:223:GLN:HE21	1:A:281:GLN:HE22	1.13	0.94
1:A:305:LYS:HE2	1:A:331:LEU:HD22	1.47	0.94
1:B:89:ARG:HH11	1:B:90:ARG:HH21	1.19	0.86
1:A:336:HIS:O	1:A:339:GLU:HG2	1.75	0.86
2:A:1201:B71:OAD	3:A:1388:HOH:O	1.93	0.85
1:B:52:ASN:HD21	1:B:57:LEU:H	1.25	0.83
1:A:38:LYS:HD3	1:A:89:ARG:HG3	1.58	0.82
1:B:55:MET:HE3	1:B:190:SER:HB2	1.60	0.82
1:B:55:MET:CE	1:B:187:GLU:HA	2.11	0.81
1:B:330:LEU:HG	3:B:1352:HOH:O	1.82	0.78
1:A:85:ASN:ND2	3:A:1446:HOH:O	2.15	0.76
1:B:172:MET:HE1	1:B:204:ASN:HB3	1.69	0.75
1:B:120:SER:HA	1:B:129:TYR:HE1	1.51	0.74
1:B:258:GLN:HG2	1:B:286:ASP:OD2	1.87	0.74
1:B:120:SER:HA	1:B:129:TYR:CE1	2.24	0.73
2:A:1201:B71:OAG	3:A:1388:HOH:O	2.05	0.73
1:A:180:ARG:HH11	1:A:204:ASN:HD21	1.34	0.72
1:B:180:ARG:HH11	1:B:204:ASN:HD21	1.39	0.69
1:A:180:ARG:HH11	1:A:204:ASN:ND2	1.90	0.68
2:B:1202:B71:OAD	3:B:1456:HOH:O	2.11	0.68
1:A:223:GLN:NE2	1:A:281:GLN:HE22	1.90	0.68
2:B:1202:B71:OAG	3:B:1456:HOH:O	2.13	0.65
1:A:116:MET:HE2	1:B:116:MET:SD	2.38	0.64
1:A:281:GLN:HA	1:A:281:GLN:HE21	1.62	0.64
1:B:172:MET:HE2	3:B:1254:HOH:O	1.97	0.64
1:A:96:HIS:HE1	3:A:1446:HOH:O	1.81	0.64
1:A:35:ILE:HG13	1:A:38:LYS:HD2	1.79	0.63
1:B:336:HIS:O	1:B:339:GLU:HG2	1.98	0.63
1:B:89:ARG:HH11	1:B:90:ARG:NH2	1.92	0.63
1:B:338:SER:HB3	3:B:1457:HOH:O	1.98	0.63
1:A:258:GLN:HE21	1:A:258:GLN:N	1.91	0.63
1:B:172:MET:CE	1:B:208:ILE:HD11	2.27	0.62
1:B:180:ARG:HH11	1:B:204:ASN:ND2	1.98	0.61
1:B:330:LEU:HD13	3:B:1357:HOH:O	1.99	0.61
1:B:55:MET:HE2	1:B:187:GLU:CA	2.22	0.61
1:A:140:LEU:HD12	1:B:113:PHE:CE2	2.36	0.60
1:A:213:ARG:NH2	3:A:1426:HOH:O	2.36	0.59
1:B:258:GLN:HB2	3:B:1355:HOH:O	2.01	0.58
1:A:223:GLN:HE21	1:A:281:GLN:NE2	1.93	0.58
1:B:203:ILE:HD12	1:B:206:LEU:HD12	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:ASN:OD1	3:A:1478:HOH:O	2.17	0.57
1:A:144:HIS:HE1	3:B:1402:HOH:O	1.85	0.57
1:A:140:LEU:O	1:A:144:HIS:HD2	1.87	0.57
1:A:152:TYR:OH	1:B:30:LYS:NZ	2.36	0.57
1:A:52:ASN:ND2	1:A:57:LEU:H	1.92	0.57
1:A:125:LYS:HG3	1:A:127:PRO:HD2	1.85	0.57
1:A:20:TRP:CH2	1:A:24:ASN:HB3	2.40	0.56
1:A:281:GLN:HA	1:A:281:GLN:NE2	2.21	0.55
1:B:47:LEU:HD13	1:B:333:ILE:HD13	1.87	0.55
1:B:45:LEU:O	1:B:49:VAL:HG23	2.07	0.54
1:B:147:GLN:HE21	1:B:147:GLN:HA	1.73	0.54
1:B:183:LEU:O	1:B:183:LEU:HD13	2.08	0.53
1:A:305:LYS:HE2	1:A:331:LEU:CD2	2.29	0.53
1:A:38:LYS:HE3	1:A:89:ARG:O	2.09	0.53
1:A:304:ILE:C	1:A:306:ASN:H	2.16	0.52
1:B:55:MET:HE3	1:B:190:SER:CB	2.35	0.52
1:A:304:ILE:C	1:A:306:ASN:N	2.68	0.52
1:B:308:ASN:C	1:B:310:ASN:H	2.18	0.52
1:A:6:MET:HE3	1:B:296:PHE:CE2	2.45	0.51
1:A:80:ASP:CB	3:A:1401:HOH:O	2.34	0.51
1:A:11:ASP:OD2	1:B:311:LYS:NZ	2.43	0.51
1:A:299:GLN:O	1:A:303:MET:HG3	2.11	0.51
1:A:183:LEU:HD22	1:A:203:ILE:HB	1.94	0.50
1:A:38:LYS:CD	1:A:89:ARG:HG3	2.35	0.49
1:B:183:LEU:HD13	1:B:183:LEU:C	2.36	0.49
1:A:329:GLU:N	3:A:1470:HOH:O	2.45	0.49
1:B:260:ASN:O	1:B:264:ARG:HG3	2.12	0.48
1:A:125:LYS:CG	1:A:127:PRO:HD2	2.44	0.48
1:A:304:ILE:O	1:A:306:ASN:N	2.47	0.48
1:B:304:ILE:HG23	1:B:313:LEU:HD21	1.95	0.48
2:B:1202:B71:HAK	2:B:1202:B71:HAT1	1.69	0.48
1:A:213:ARG:NH1	1:A:334:ILE:O	2.46	0.48
1:A:197:HIS:HD2	1:A:198:SER:O	1.98	0.47
1:A:21:SER:H	1:A:24:ASN:HD22	1.63	0.47
1:B:20:TRP:CH2	1:B:24:ASN:HB3	2.50	0.47
1:A:305:LYS:HG2	1:A:331:LEU:HD21	1.97	0.46
1:B:43:PHE:CD1	1:B:339:GLU:HG3	2.50	0.46
1:A:230:PHE:CD1	1:A:230:PHE:N	2.83	0.46
1:B:52:ASN:ND2	1:B:57:LEU:H	2.03	0.46
1:B:340:LEU:HB3	2:B:1202:B71:HAQ2	1.98	0.45
1:A:141:ILE:HG23	1:B:110:TYR:HE1	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:303:MET:HE3	1:B:7:GLU:HB3	1.99	0.45
1:B:79:ILE:HD12	1:B:112:TYR:OH	2.17	0.45
1:B:75:SER:OG	1:B:111:MET:HB2	2.18	0.44
1:B:178:LEU:HB2	3:B:1238:HOH:O	2.16	0.44
1:A:96:HIS:HD2	3:A:1384:HOH:O	2.00	0.44
1:B:178:LEU:HD13	2:B:1202:B71:HAN1	1.99	0.44
1:B:251:LYS:HB2	1:B:251:LYS:NZ	2.32	0.44
1:A:90:ARG:NH2	3:A:1381:HOH:O	2.51	0.44
1:A:47:LEU:O	1:A:51:ILE:HG12	2.18	0.43
1:B:183:LEU:HD13	1:B:187:GLU:HG3	2.00	0.43
1:A:241:PHE:HB3	1:A:242:PRO:HD3	2.00	0.43
1:A:125:LYS:NZ	1:A:127:PRO:HB2	2.32	0.43
1:B:21:SER:H	1:B:24:ASN:HD22	1.65	0.43
1:A:180:ARG:HD3	1:A:204:ASN:HD21	1.82	0.43
1:A:255:GLN:HB3	1:A:258:GLN:NE2	2.34	0.43
1:B:331:LEU:HD23	1:B:335:ASP:OD2	2.19	0.43
1:A:144:HIS:CE1	3:B:1402:HOH:O	2.67	0.43
1:B:214:ASP:OD2	3:B:1456:HOH:O	2.20	0.42
1:A:281:GLN:HE21	1:A:281:GLN:CA	2.31	0.42
1:B:251:LYS:NZ	1:B:251:LYS:CB	2.82	0.42
1:B:49:VAL:O	1:B:50:GLN:C	2.62	0.42
1:B:119:VAL:HG11	1:B:133:ILE:HD13	2.02	0.42
1:B:286:ASP:OD1	1:B:286:ASP:C	2.61	0.42
1:A:114:ARG:HA	1:A:114:ARG:HD2	1.87	0.41
1:A:21:SER:H	1:A:24:ASN:ND2	2.18	0.41
1:B:278:LYS:O	1:B:281:GLN:HG2	2.20	0.41
1:B:167:TYR:CD2	1:B:167:TYR:C	2.98	0.41
1:A:113:PHE:CE2	1:B:140:LEU:HD13	2.56	0.41
1:B:101:VAL:HB	1:B:102:PRO:HD3	2.02	0.41
1:B:180:ARG:O	1:B:184:ARG:HG3	2.21	0.41
1:B:295:ASN:O	1:B:299:GLN:HG3	2.21	0.41
1:A:47:LEU:HD13	1:A:333:ILE:HD13	2.03	0.41
1:A:90:ARG:CZ	3:A:1381:HOH:O	2.69	0.41
1:A:210:TYR:CD2	1:A:210:TYR:C	2.99	0.41
1:B:152:TYR:CD2	1:B:152:TYR:C	2.98	0.41
1:B:340:LEU:HD23	1:B:340:LEU:HA	1.86	0.41
1:A:38:LYS:HE2	1:A:89:ARG:HB3	2.04	0.40
1:A:140:LEU:CD1	1:B:113:PHE:CE2	3.03	0.40
1:A:170:MET:SD	1:A:174:LYS:HE2	2.61	0.40

All (1) 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)
3:B:1266:HOH:O	3:B:1460:HOH:O[3_655]	2.18	0.02

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	295/340 (87%)	293 (99%)	1 (0%)	1 (0%)	36	35
1	B	298/340 (88%)	291 (98%)	7 (2%)	0	100	100
All	All	593/680 (87%)	584 (98%)	8 (1%)	1 (0%)	43	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	305	LYS

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	285/317 (90%)	275 (96%)	10 (4%)	32	32
1	B	290/317 (92%)	282 (97%)	8 (3%)	38	41
All	All	575/634 (91%)	557 (97%)	18 (3%)	35	37

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

Mol	Chain	Res	Type
1	A	6	MET
1	A	27	LEU
1	A	61	GLN
1	A	158	PRO
1	A	178	LEU
1	A	203	ILE
1	A	230	PHE
1	A	258	GLN
1	A	277	LEU
1	A	339	GLU
1	B	90	ARG
1	B	120	SER
1	B	147	GLN
1	B	178	LEU
1	B	181	LEU
1	B	251	LYS
1	B	258	GLN
1	B	339	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	52	ASN
1	A	61	GLN
1	A	73	HIS
1	A	85	ASN
1	A	96	HIS
1	A	106	ASN
1	A	109	ASN
1	A	130	HIS
1	A	144	HIS
1	A	197	HIS
1	A	204	ASN
1	A	255	GLN
1	A	258	GLN
1	A	259	HIS
1	A	260	ASN
1	A	281	GLN
1	A	298	ASN
1	A	306	ASN
1	B	52	ASN
1	B	109	ASN

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Mol	Chain	Res	Type
1	B	121	GLN
1	B	137	ASN
1	B	147	GLN
1	B	204	ASN
1	B	255	GLN
1	B	281	GLN
1	B	288	ASN
1	B	298	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](#)

2 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	$\# Z > 2$
2	B71	A	1201	-	26,27,27	2.72	8 (30%)	31,36,36	1.20	4 (12%)
2	B71	B	1202	-	26,27,27	2.76	8 (30%)	31,36,36	1.08	3 (9%)

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
2	B71	A	1201	-	-	1/25/27/27	0/1/1/1
2	B71	B	1202	-	-	1/25/27/27	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1202	B71	PAZ-OAE	7.19	1.61	1.49
2	B	1202	B71	PBA-OAF	7.15	1.61	1.49
2	A	1201	B71	PAZ-OAE	7.13	1.61	1.49
2	A	1201	B71	PBA-OAF	6.89	1.61	1.49
2	B	1202	B71	PAZ-OAD	4.21	1.61	1.54
2	B	1202	B71	PBA-OAG	4.18	1.61	1.54
2	B	1202	B71	PBA-OAC	-4.11	1.48	1.54
2	A	1201	B71	PBA-OAG	4.11	1.61	1.54
2	A	1201	B71	PBA-OAC	-4.09	1.48	1.54
2	A	1201	B71	PAZ-OAD	4.04	1.61	1.54
2	A	1201	B71	PAZ-OAB	-4.03	1.48	1.54
2	B	1202	B71	PAZ-OAB	-3.99	1.48	1.54
2	B	1202	B71	CAK-NAY	3.51	1.40	1.35
2	A	1201	B71	CAK-NAY	3.49	1.40	1.35
2	A	1201	B71	CAJ-NAY	2.51	1.40	1.34
2	B	1202	B71	CAJ-NAY	2.46	1.40	1.34

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1201	B71	CAT-OAV-CAW	-3.22	109.56	117.93
2	A	1201	B71	OAD-PAZ-OAE	-2.73	106.61	113.45
2	B	1202	B71	OAD-PAZ-OAE	-2.62	106.87	113.45
2	B	1202	B71	OAG-PBA-OAF	-2.54	107.07	113.45
2	A	1201	B71	OAG-PBA-OAF	-2.31	107.67	113.45
2	A	1201	B71	OAF-PBA-CAX	-2.29	105.79	112.29
2	B	1202	B71	OAF-PBA-CAX	-2.19	106.07	112.29

There are no chirality outliers.

All (2) torsion outliers are listed below:

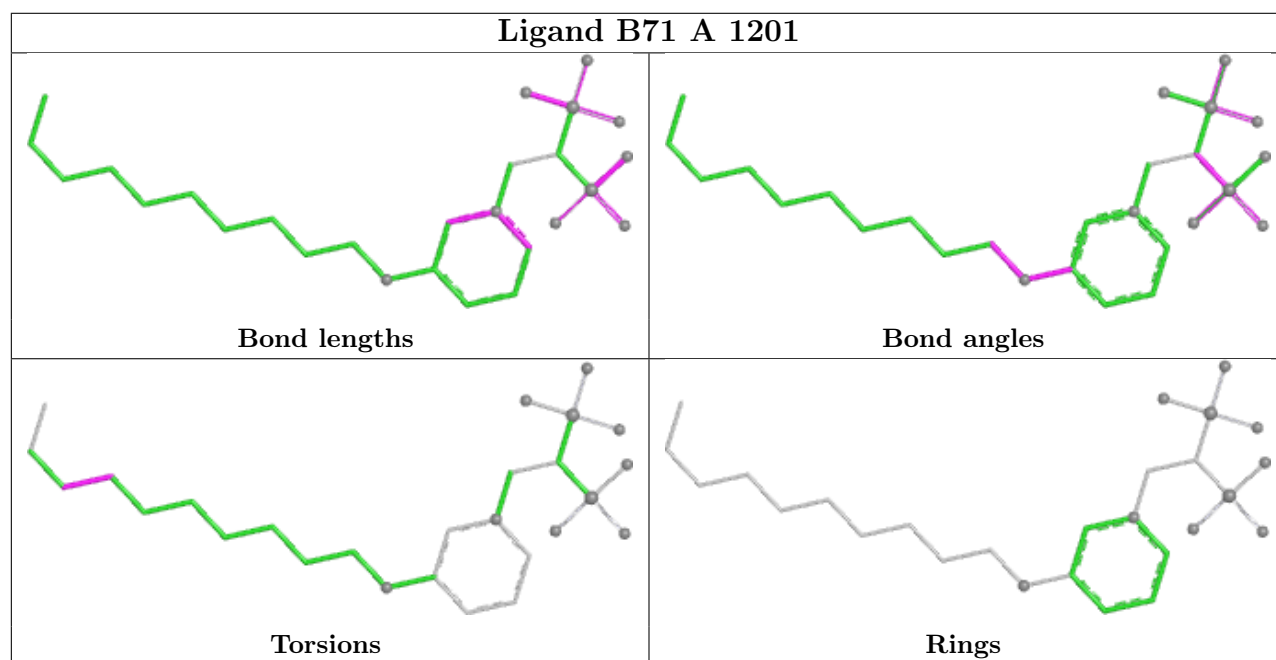
Mol	Chain	Res	Type	Atoms
2	A	1201	B71	CAL-CAM-CAN-CAO
2	B	1202	B71	CAK-CAW-OAV-CAT

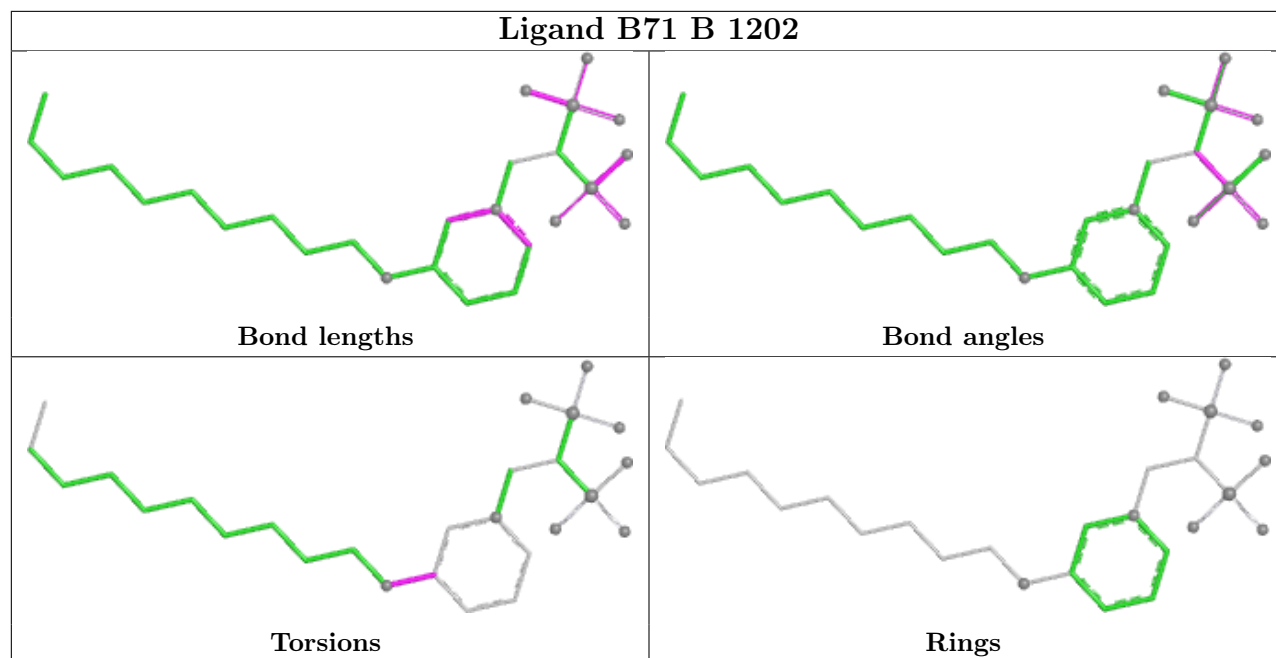
There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1201	B71	2	0
2	B	1202	B71	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	305/340 (89%)	-0.07	8 (2%) 57 56	26, 38, 68, 81	0
1	B	308/340 (90%)	0.18	16 (5%) 33 31	26, 42, 79, 94	0
All	All	613/680 (90%)	0.06	24 (3%) 43 42	26, 40, 72, 94	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	331	LEU	4.1
1	B	314	PRO	3.3
1	A	331	LEU	3.3
1	B	8	ALA	3.2
1	B	37	LEU	3.1
1	B	191	PRO	3.0
1	A	38	LYS	3.0
1	B	332	TYR	2.6
1	B	313	LEU	2.6
1	B	340	LEU	2.6
1	A	230	PHE	2.5
1	B	339	GLU	2.4
1	B	335	ASP	2.3
1	B	4	ASN	2.3
1	B	42	ASN	2.3
1	B	330	LEU	2.3
1	B	192	SER	2.3
1	B	333	ILE	2.3
1	A	197	HIS	2.2
1	A	60	ASP	2.1
1	A	339	GLU	2.1
1	A	229	GLY	2.1
1	B	60	ASP	2.0
1	A	335	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](#)

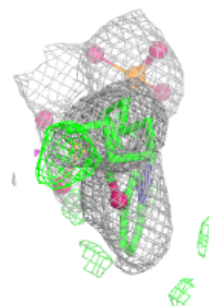
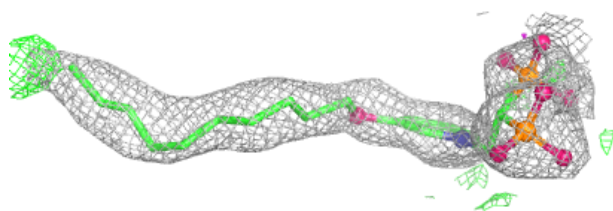
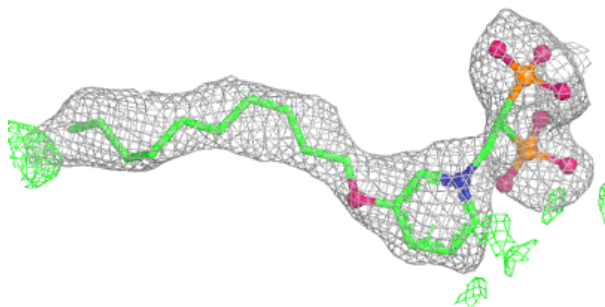
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	B71	B	1202	27/27	0.84	0.14	72,75,78,78	0
2	B71	A	1201	27/27	0.88	0.11	55,62,64,66	0

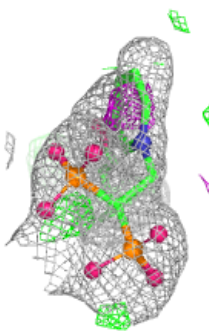
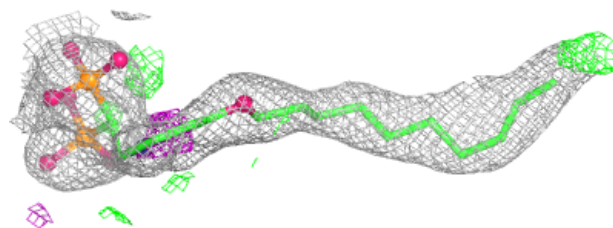
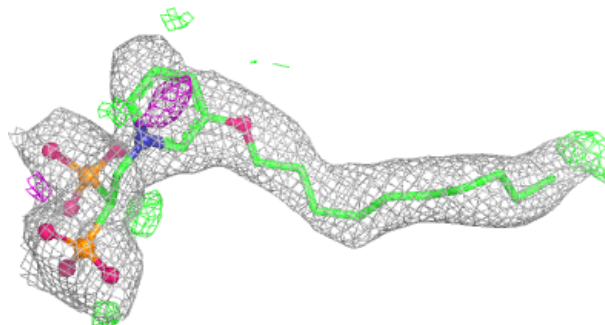
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around B71 B 1202:

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

**Electron density around B71 A 1201:**

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



6.5 Other polymers ⓘ

There are no such residues in this entry.