



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 12:46 AM UTC

PDB ID : 7E40 / pdb_00007e40
Title : Mechanism of Phosphate Sensing and Signaling Revealed by Rice SPX1-PHR2 Complex Structure
Authors : Zhou, J.; Hu, Q.; Yao, D.; Xing, W.
Deposited on : 2021-02-09
Resolution : 2.60 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

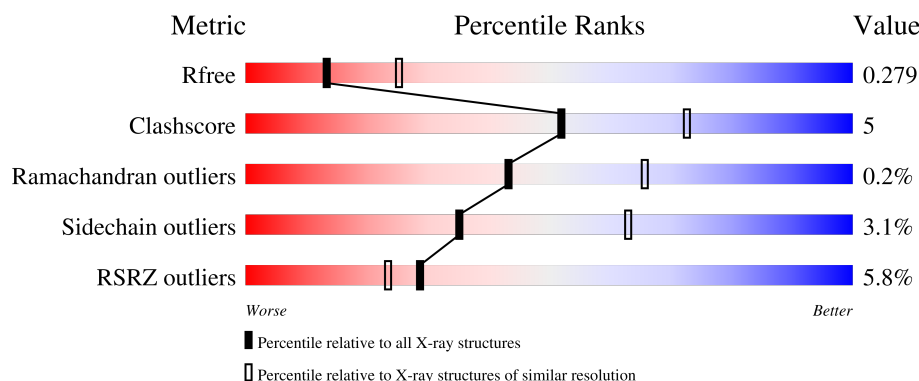
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	134	 5% 72% 6% • 21%
1	C	134	 5% 65% 10% • 24%
2	B	358	 2% 83% 9% • 7%
2	D	358	 9% 81% 14% • •

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7318 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 Protein PHOSPHATE STARVATION RESPONSE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	106	Total	C	N	O	S	0	0	0
			880	549	166	156	9			
1	C	102	Total	C	N	O	S	0	0	0
			848	531	157	151	9			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	247	SER	-	expression tag	UNP Q6Z156
A	263	MET	VAL	engineered mutation	UNP Q6Z156
A	278	MET	LEU	engineered mutation	UNP Q6Z156
A	295	MET	LEU	engineered mutation	UNP Q6Z156
A	340	MET	LEU	engineered mutation	UNP Q6Z156
C	247	SER	-	expression tag	UNP Q6Z156
C	263	MET	VAL	engineered mutation	UNP Q6Z156
C	278	MET	LEU	engineered mutation	UNP Q6Z156
C	295	MET	LEU	engineered mutation	UNP Q6Z156
C	340	MET	LEU	engineered mutation	UNP Q6Z156

- Molecule 2 is a protein called SPX domain-containing protein 1,Endolysin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	332	Total	C	N	O	S	0	0	0
			2694	1719	470	494	11			
2	D	345	Total	C	N	O	S	0	0	0
			2771	1755	491	515	10			

There are 6 discrepancies between the modelled and reference sequences:

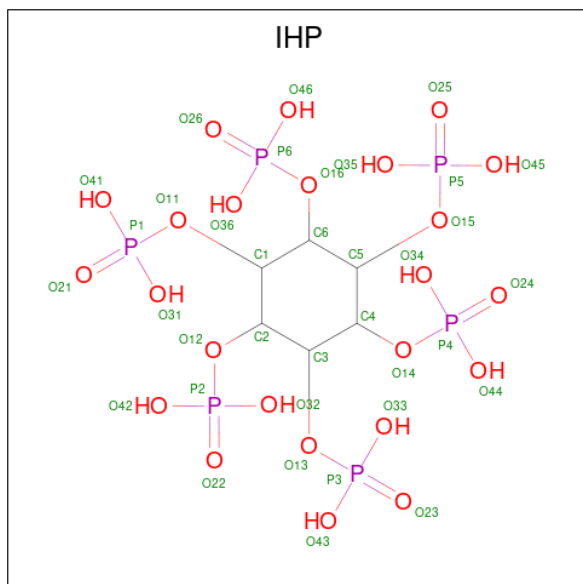
Chain	Residue	Modelled	Actual	Comment	Reference
B	0	SER	-	expression tag	UNP Q69XJ0
B	250	THR	CYS	engineered mutation	UNP D9IEF7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	293	ALA	CYS	engineered mutation	UNP D9IEF7
D	0	SER	-	expression tag	UNP Q69XJ0
D	250	THR	CYS	engineered mutation	UNP D9IEF7
D	293	ALA	CYS	engineered mutation	UNP D9IEF7

- Molecule 3 is INOSITOL HEXAKISPHOSPHATE (CCD ID: IHP) (formula: $C_6H_{18}O_{24}P_6$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	O	P	0	0
			36	6	24	6		
3	B	1	Total	C	O	P	0	0
			36	6	24	6		
3	D	1	Total	C	O	P	0	0
			36	6	24	6		

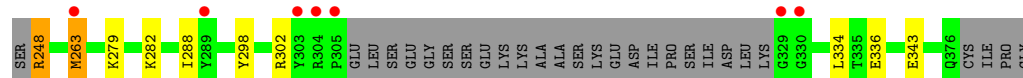
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	5	Total	O	0	0
			5	5		
4	B	9	Total	O	0	0
			9	9		
4	D	3	Total	O	0	0
			3	3		

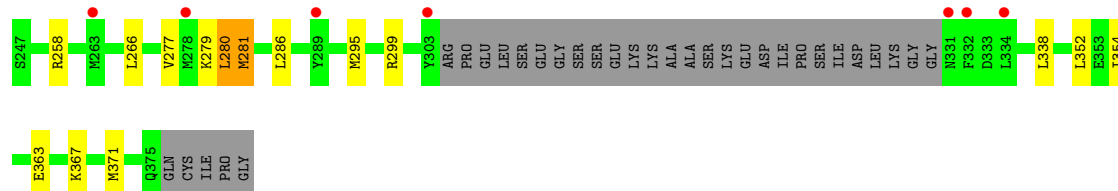
3 Residue-property plots [i](#)

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.

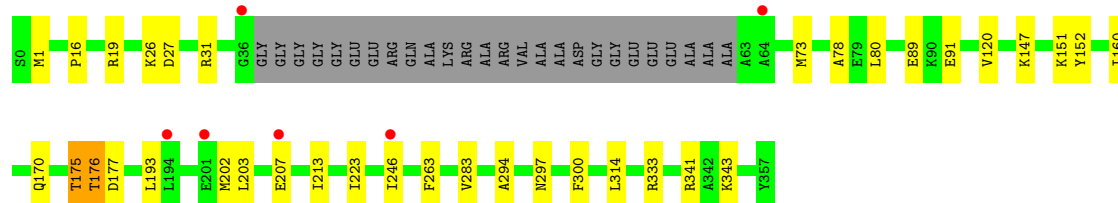
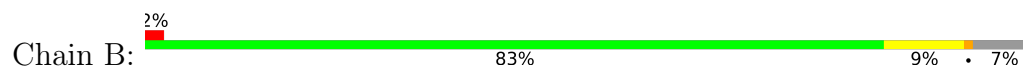
- Molecule 1: Protein PHOSPHATE STARVATION RESPONSE 2



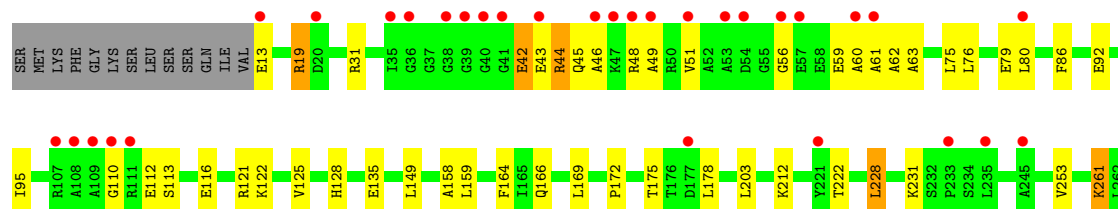
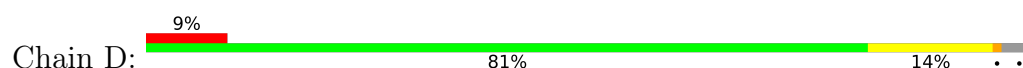
- Molecule 1: Protein PHOSPHATE STARVATION RESPONSE 2



- Molecule 2: SPX domain-containing protein 1,Endolysin



- Molecule 2: SPX domain-containing protein 1,Endolysin



F263	
F300	
R321	
K343	
W354	
Y357	

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.24Å 107.49Å 174.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.76 – 2.60 45.76 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (45.76-2.60) 99.9 (45.76-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.70 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.18_3861	Depositor
R, R_{free}	0.226 , 0.274 0.230 , 0.279	Depositor DCC
R_{free} test set	1935 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	53.6	Xtriage
Anisotropy	0.297	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.2	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	0.93	EDS
Total number of atoms	7318	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.69% 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: IHP

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	0.79	0/892	1.47	3/1188 (0.3%)
1	C	0.79	1/859 (0.1%)	1.49	2/1144 (0.2%)
2	B	0.81	0/2737	1.45	7/3675 (0.2%)
2	D	0.80	0/2814	1.47	18/3779 (0.5%)
All	All	0.80	1/7302 (0.0%)	1.47	30/9786 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	281	MET	SD-CE	-5.49	1.65	1.79

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	44	ARG	N-CA-C	-7.42	102.81	112.68
2	D	128	HIS	CA-C-N	5.86	126.48	119.98
2	D	128	HIS	C-N-CA	5.86	126.48	119.98
2	D	86	PHE	CA-C-N	5.84	128.03	120.44
2	D	86	PHE	C-N-CA	5.84	128.03	120.44
2	B	78	ALA	CA-C-N	5.78	127.95	120.44
2	B	78	ALA	C-N-CA	5.78	127.95	120.44
2	B	175	THR	O-C-N	5.67	128.86	122.22
1	A	288	ILE	CA-C-N	5.38	127.44	120.44
1	A	288	ILE	C-N-CA	5.38	127.44	120.44
2	B	333	ARG	CA-C-N	5.37	127.74	120.38
2	B	333	ARG	C-N-CA	5.37	127.74	120.38
2	D	56	GLY	CA-C-O	-5.31	115.92	121.28
2	D	110	GLY	CA-C-N	5.30	129.83	121.56
2	D	110	GLY	C-N-CA	5.30	129.83	121.56
1	C	258	ARG	CA-C-N	5.24	127.25	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	258	ARG	C-N-CA	5.24	127.25	120.44
2	D	231	LYS	CA-C-N	5.23	127.66	120.39
2	D	231	LYS	C-N-CA	5.23	127.66	120.39
2	D	354	TRP	CA-C-N	5.19	127.49	120.38
2	D	354	TRP	C-N-CA	5.19	127.49	120.38
2	D	321	ARG	CA-C-N	5.16	127.14	120.44
2	D	321	ARG	C-N-CA	5.16	127.14	120.44
2	D	92	GLU	CA-C-N	5.14	127.13	120.44
2	D	92	GLU	C-N-CA	5.14	127.13	120.44
1	A	343	GLU	CB-CG-CD	5.12	121.31	112.60
2	D	79	GLU	CA-C-N	5.05	127.01	120.44
2	D	79	GLU	C-N-CA	5.05	127.01	120.44
2	B	151	LYS	CA-C-N	5.03	126.98	120.44
2	B	151	LYS	C-N-CA	5.03	126.98	120.44

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	880	0	901	4	0
1	C	848	0	868	8	0
2	B	2694	0	2771	20	0
2	D	2771	0	2822	48	0
3	B	72	0	12	0	0
3	D	36	0	6	0	0
4	A	5	0	0	0	0
4	B	9	0	0	0	0
4	D	3	0	0	0	0
All	All	7318	0	7380	75	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 5.

All (75) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:51:VAL:HG11	2:D:159:LEU:HG	1.46	0.97
2:D:112:GLU:HG2	2:D:113:SER:H	1.36	0.87
2:D:112:GLU:OE1	2:D:116:GLU:HG2	1.77	0.83
2:D:76:LEU:O	2:D:80:LEU:HD13	1.79	0.82
2:D:51:VAL:CG1	2:D:159:LEU:HG	2.11	0.79
1:C:277:VAL:HG12	1:C:281:MET:HE2	1.65	0.79
2:D:51:VAL:HG12	2:D:159:LEU:HB2	1.66	0.76
2:D:51:VAL:HG11	2:D:159:LEU:CG	2.14	0.76
2:B:203:LEU:HD21	2:B:297:ASN:ND2	2.01	0.75
2:D:59:GLU:OE2	2:D:63:ALA:HA	1.85	0.74
2:D:112:GLU:HG2	2:D:113:SER:N	2.03	0.72
2:B:91:GLU:OE2	2:B:177:ASP:HB3	1.89	0.70
2:D:60:ALA:O	2:D:62:ALA:N	2.24	0.70
1:A:263:MET:HE1	1:A:298:TYR:CD2	2.29	0.68
2:D:222:THR:HG22	2:D:228:LEU:HA	1.74	0.68
2:D:51:VAL:HG13	2:D:159:LEU:N	2.09	0.67
2:B:203:LEU:CD2	2:B:297:ASN:ND2	2.58	0.67
2:D:51:VAL:CG1	2:D:159:LEU:CG	2.74	0.66
2:D:203:LEU:HD12	2:D:263:PHE:CZ	2.34	0.63
2:D:203:LEU:HD12	2:D:263:PHE:HZ	1.64	0.62
2:D:112:GLU:CG	2:D:113:SER:H	2.08	0.62
2:D:45:GLN:O	2:D:48:ARG:HG2	2.00	0.62
2:D:80:LEU:HD21	2:D:164:PHE:CE1	2.37	0.59
2:D:80:LEU:HD21	2:D:164:PHE:CZ	2.37	0.59
2:B:175:THR:O	2:B:176:THR:HB	2.02	0.59
2:D:212:LYS:HG2	2:D:253:VAL:HG22	1.85	0.57
2:D:51:VAL:HG12	2:D:159:LEU:CB	2.33	0.57
2:D:51:VAL:CG1	2:D:159:LEU:N	2.69	0.56
2:B:283:VAL:HG11	2:B:314:LEU:HD22	1.88	0.56
2:D:76:LEU:O	2:D:80:LEU:CD1	2.52	0.56
2:B:203:LEU:HD13	2:B:300:PHE:HD2	1.72	0.55
2:B:207:GLU:HG2	2:B:341:ARG:HH12	1.72	0.55
2:D:95:ILE:HD11	2:D:178:LEU:HD21	1.90	0.54
2:D:51:VAL:CG1	2:D:159:LEU:CB	2.86	0.54
2:B:203:LEU:CD1	2:B:300:PHE:HD2	2.21	0.54
2:B:203:LEU:HD13	2:B:300:PHE:CD2	2.44	0.53
2:B:1:MET:HE2	1:C:354:ILE:HG23	1.91	0.53
2:D:59:GLU:CG	2:D:63:ALA:HA	2.40	0.52
1:A:248:ARG:CZ	2:B:89:GLU:HG3	2.40	0.52
2:D:203:LEU:HD13	2:D:300:PHE:CD2	2.46	0.51
1:C:338:LEU:HD21	2:D:122:LYS:HG3	1.92	0.51
2:D:80:LEU:CD2	2:D:164:PHE:CZ	2.95	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:203:LEU:HD21	2:B:297:ASN:CG	2.35	0.50
2:B:120:VAL:HG13	2:B:193:LEU:HD21	1.94	0.49
2:B:203:LEU:CD1	2:B:300:PHE:CD2	2.96	0.49
2:D:172:PRO:HA	2:D:175:THR:HG22	1.94	0.48
1:C:266:LEU:HD11	1:C:280:LEU:HD23	1.95	0.48
2:D:45:GLN:O	2:D:48:ARG:NE	2.45	0.48
2:B:27:ASP:O	2:B:31:ARG:HG2	2.15	0.47
2:D:203:LEU:CD1	2:D:300:PHE:CD2	2.98	0.47
1:C:352:LEU:HD13	2:D:135:GLU:HG3	1.96	0.46
1:C:295:MET:HG3	1:C:299:ARG:HD2	1.98	0.46
2:D:51:VAL:CG1	2:D:159:LEU:HB2	2.42	0.46
2:D:51:VAL:HG11	2:D:159:LEU:CD1	2.46	0.45
1:C:367:LYS:HG2	1:C:371:MET:HE2	1.98	0.45
2:D:42:GLU:C	2:D:44:ARG:H	2.25	0.44
2:D:51:VAL:HG13	2:D:158:ALA:C	2.43	0.43
2:D:45:GLN:O	2:D:48:ARG:CG	2.65	0.43
2:D:112:GLU:CG	2:D:113:SER:N	2.71	0.43
2:D:149:LEU:HD11	2:D:164:PHE:CD1	2.53	0.43
2:D:261:LYS:HZ3	2:D:261:LYS:HA	1.84	0.42
2:D:121:ARG:O	2:D:125:VAL:HG23	2.19	0.42
2:D:203:LEU:CD1	2:D:300:PHE:HD2	2.33	0.42
1:C:363:GLU:HG3	2:D:169:LEU:HD11	2.02	0.42
2:B:213:ILE:HG13	2:B:223:ILE:HD12	2.00	0.42
2:D:43:GLU:HA	2:D:46:ALA:HB3	2.02	0.41
2:B:73:MET:HE1	2:B:152:TYR:CZ	2.54	0.41
2:B:202:MET:HE1	2:B:294:ALA:HA	2.02	0.41
1:A:279:LYS:O	1:A:282:LYS:HG2	2.21	0.41
2:B:16:PRO:O	2:B:19:ARG:CB	2.69	0.41
2:D:59:GLU:CD	2:D:63:ALA:HA	2.43	0.41
2:D:112:GLU:CD	2:D:116:GLU:HG2	2.43	0.41
2:D:48:ARG:HG3	2:D:49:ALA:N	2.36	0.41
1:A:298:TYR:CZ	1:A:302:ARG:HD2	2.56	0.41
2:B:203:LEU:HD12	2:B:263:PHE:CZ	2.56	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	102/134 (76%)	102 (100%)	0	0	100	100
1	C	98/134 (73%)	97 (99%)	1 (1%)	0	100	100
2	B	328/358 (92%)	319 (97%)	9 (3%)	0	100	100
2	D	343/358 (96%)	326 (95%)	15 (4%)	2 (1%)	21	42
All	All	871/984 (88%)	844 (97%)	25 (3%)	2 (0%)	43	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	61	ALA
2	D	19	ARG

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	95/119 (80%)	91 (96%)	4 (4%)	26	52
1	C	92/119 (77%)	89 (97%)	3 (3%)	33	61
2	B	288/300 (96%)	280 (97%)	8 (3%)	38	66
2	D	288/300 (96%)	279 (97%)	9 (3%)	35	63
All	All	763/838 (91%)	739 (97%)	24 (3%)	35	63

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	248	ARG
1	A	263	MET
1	A	334	LEU
1	A	336	GLU
2	B	26	LYS
2	B	80	LEU
2	B	147	LYS
2	B	160	ILE
2	B	170	GLN
2	B	176	THR
2	B	246	ILE
2	B	343	LYS
1	C	279	LYS
1	C	280	LEU
1	C	286	LEU
2	D	13	GLU
2	D	19	ARG
2	D	31	ARG
2	D	42	GLU
2	D	75	LEU
2	D	166	GLN
2	D	228	LEU
2	D	261	LYS
2	D	343	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	170	GLN
2	B	251	ASN
1	C	341	GLN
1	C	345	GLN
2	D	84	ASN
2	D	98	GLN
2	D	141	ASN
2	D	166	GLN
2	D	171	GLN
2	D	185	GLN
2	D	197	ASN
2	D	264	ASN
2	D	312	ASN
2	D	328	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

3 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	IHP	B	402	-	36,36,36	1.67	6 (16%)	60,60,60	1.21	5 (8%)
3	IHP	B	401	-	36,36,36	1.70	6 (16%)	60,60,60	1.31	7 (11%)
3	IHP	D	401	-	36,36,36	1.70	6 (16%)	60,60,60	1.20	5 (8%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	IHP	B	402	-	-	3/30/54/54	0/1/1/1
3	IHP	B	401	-	-	12/30/54/54	0/1/1/1
3	IHP	D	401	-	-	2/30/54/54	0/1/1/1

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	401	IHP	P2-O12	3.92	1.66	1.59
3	B	401	IHP	P1-O11	3.88	1.66	1.59
3	B	402	IHP	P5-O15	3.88	1.66	1.59
3	D	401	IHP	P5-O15	3.80	1.66	1.59
3	B	402	IHP	P6-O16	3.80	1.66	1.59
3	D	401	IHP	P4-O14	3.72	1.66	1.59
3	B	401	IHP	P4-O14	3.70	1.66	1.59
3	D	401	IHP	P3-O13	3.67	1.66	1.59
3	D	401	IHP	P6-O16	3.61	1.65	1.59
3	B	402	IHP	P2-O12	3.59	1.65	1.59
3	B	402	IHP	P3-O13	3.52	1.65	1.59
3	D	401	IHP	P2-O12	3.49	1.65	1.59
3	B	401	IHP	P3-O13	3.44	1.65	1.59
3	B	402	IHP	P4-O14	3.32	1.65	1.59
3	B	401	IHP	P5-O15	3.24	1.65	1.59
3	D	401	IHP	P1-O11	3.23	1.65	1.59
3	B	401	IHP	P6-O16	3.20	1.65	1.59
3	B	402	IHP	P1-O11	3.05	1.64	1.59

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	401	IHP	C6-C1-C2	4.57	120.45	110.43
3	B	401	IHP	O13-C3-C2	4.14	117.56	108.76
3	D	401	IHP	C5-C6-C1	3.67	118.47	110.43
3	B	401	IHP	O14-C4-C3	3.56	116.34	108.76
3	B	402	IHP	C6-C1-C2	3.53	118.17	110.43
3	B	401	IHP	O14-C4-C5	3.30	115.77	108.76
3	D	401	IHP	C3-C2-C1	3.21	117.47	110.43
3	B	402	IHP	C4-C3-C2	3.06	117.15	110.43
3	B	402	IHP	C3-C2-C1	3.04	117.10	110.43
3	B	402	IHP	C5-C4-C3	2.88	116.74	110.43
3	B	401	IHP	O13-C3-C4	2.77	114.64	108.76
3	B	401	IHP	C5-C6-C1	2.26	115.39	110.43
3	B	401	IHP	O12-C2-C1	2.21	113.46	108.76
3	D	401	IHP	O14-C4-C5	2.20	113.43	108.76
3	B	402	IHP	O16-C6-C1	2.08	113.19	108.76
3	B	401	IHP	C4-C3-C2	2.01	114.84	110.43
3	D	401	IHP	C5-C4-C3	2.01	114.84	110.43

There are no chirality outliers.

All (17) torsion outliers are listed below:

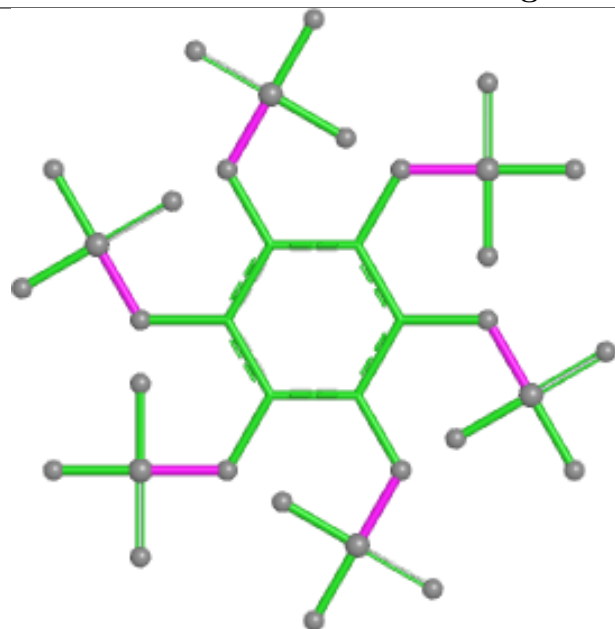
Mol	Chain	Res	Type	Atoms
3	B	401	IHP	C1-C2-O12-P2
3	B	401	IHP	C2-C3-O13-P3
3	B	401	IHP	C3-C4-O14-P4
3	B	402	IHP	C6-C5-O15-P5
3	B	401	IHP	C5-C4-O14-P4
3	B	401	IHP	C1-O11-P1-O41
3	B	401	IHP	C3-O13-P3-O33
3	D	401	IHP	C3-O13-P3-O33
3	B	401	IHP	C6-C1-O11-P1
3	B	401	IHP	C3-C2-O12-P2
3	D	401	IHP	C3-O13-P3-O23
3	B	401	IHP	C2-O12-P2-O42
3	B	402	IHP	C3-O13-P3-O33
3	B	401	IHP	C3-O13-P3-O23
3	B	401	IHP	C4-O14-P4-O24
3	B	401	IHP	C6-O16-P6-O26
3	B	402	IHP	C4-O14-P4-O24

There are no ring outliers.

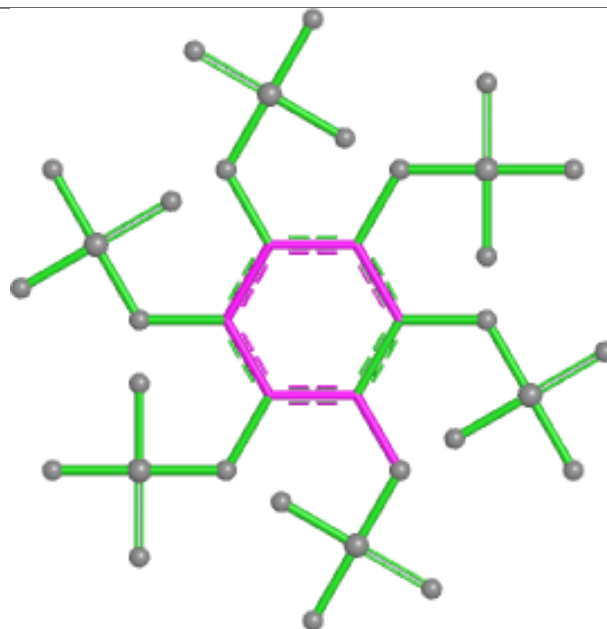
No monomer is involved in short contacts.

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.

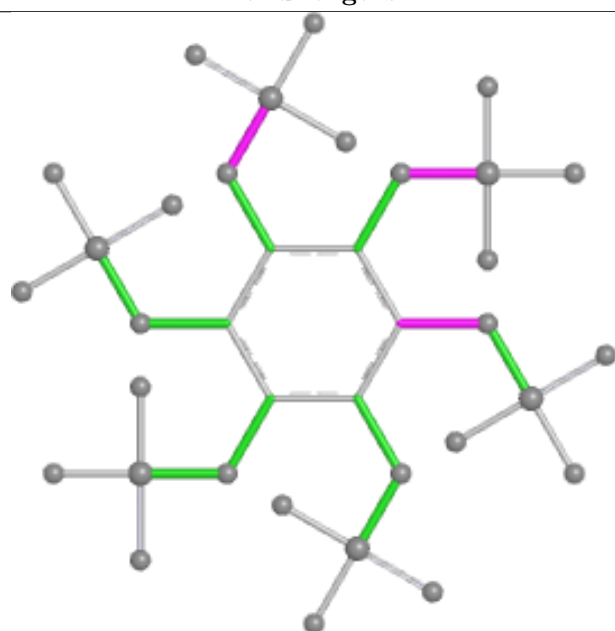
Ligand IHP B 402



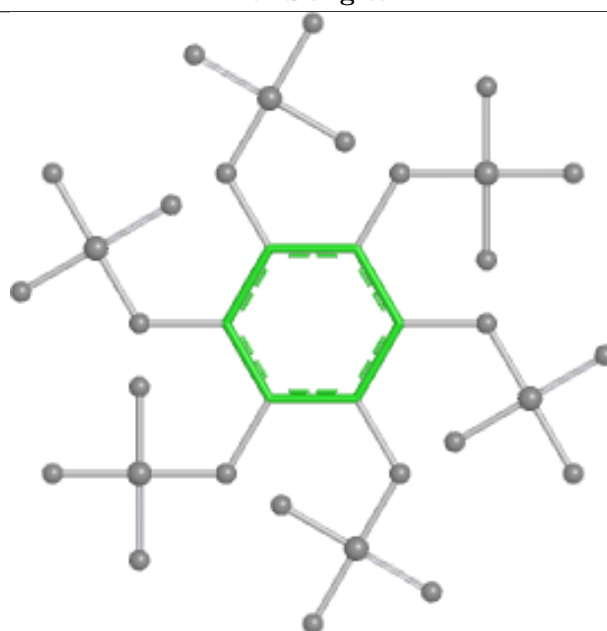
Bond lengths



Bond angles

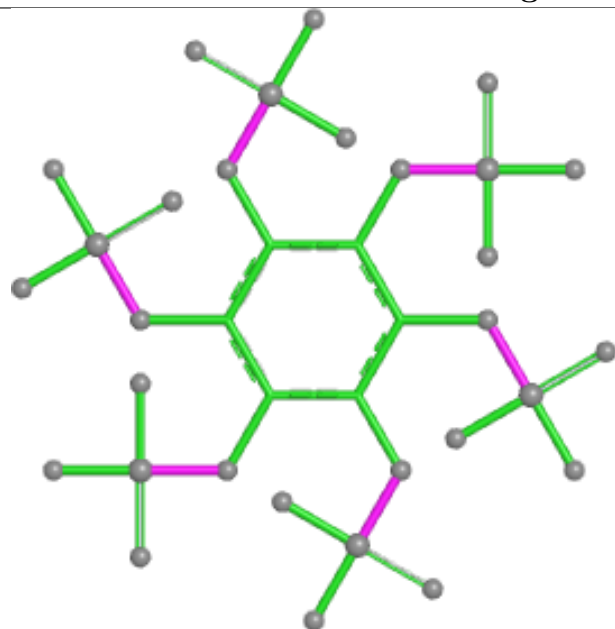


Torsions

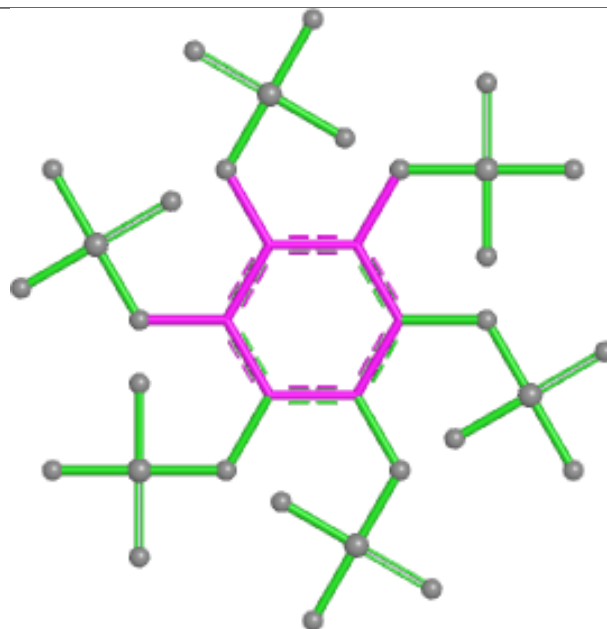


Rings

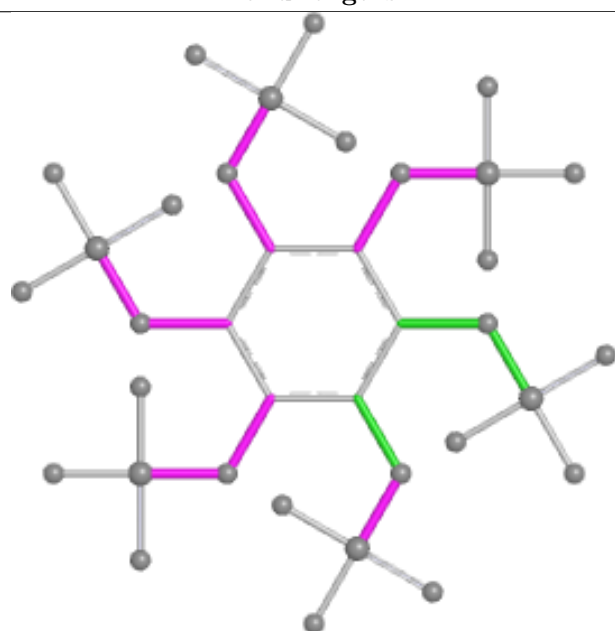
Ligand IHP B 401



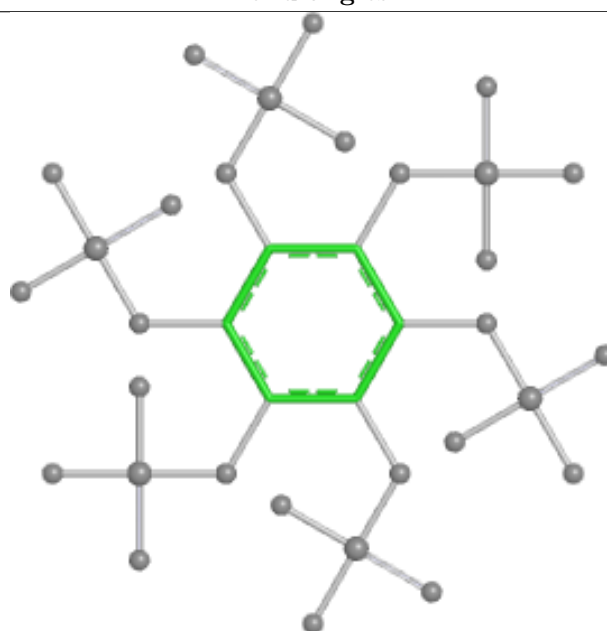
Bond lengths



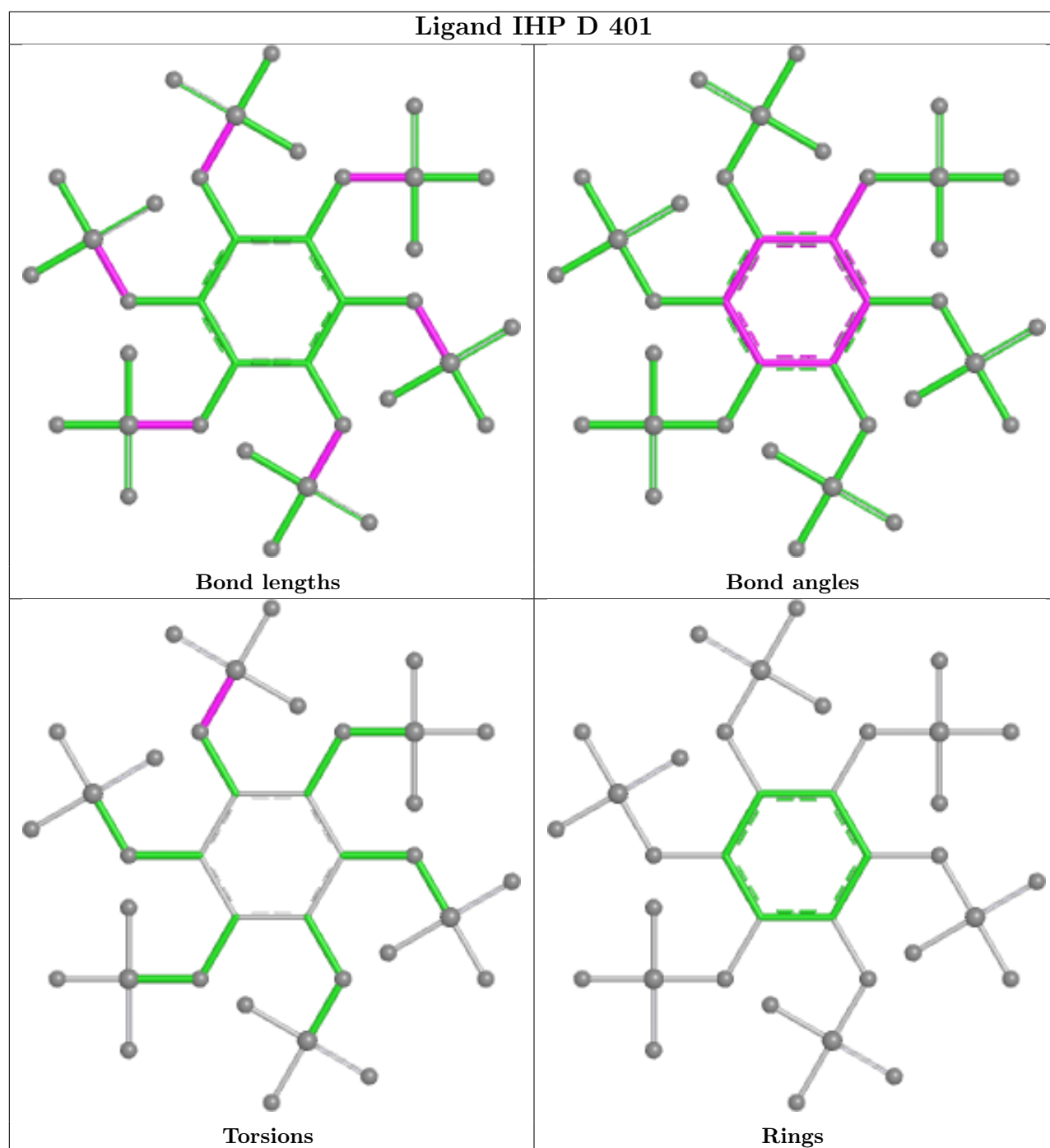
Bond angles



Torsions



Rings



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	106/134 (79%)	0.55	7 (6%)	24 19	43, 63, 103, 159	0
1	C	102/134 (76%)	0.46	7 (6%)	23 18	40, 63, 87, 128	0
2	B	332/358 (92%)	0.36	6 (1%)	67 63	41, 61, 106, 120	0
2	D	345/358 (96%)	0.52	31 (8%)	15 11	39, 62, 106, 134	0
All	All	885/984 (89%)	0.46	51 (5%)	29 23	39, 62, 106, 159	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	41	GLY	5.3
1	A	263	MET	4.5
2	D	110	GLY	4.5
2	D	40	GLY	4.3
1	A	305	PRO	3.8
2	D	60	ALA	3.8
2	D	51	VAL	3.8
1	A	329	GLY	3.8
2	D	20	ASP	3.7
1	C	303	TYR	3.7
1	C	331	ASN	3.1
2	D	39	GLY	3.1
2	B	207	GLU	3.1
2	D	107	ARG	3.1
2	D	43	GLU	3.0
2	B	36	GLY	3.0
1	A	330	GLY	2.9
2	D	36	GLY	2.8
2	D	46	ALA	2.8
2	D	57	GLU	2.8
2	D	80	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	278	MET	2.7
2	D	61	ALA	2.7
1	A	303	TYR	2.6
2	D	233	PRO	2.6
2	D	53	ALA	2.6
2	D	245	ALA	2.5
2	D	49	ALA	2.5
1	A	304	ARG	2.5
2	D	48	ARG	2.5
1	C	334	LEU	2.4
2	B	194	LEU	2.4
2	D	108	ALA	2.4
2	D	109	ALA	2.4
1	A	289	TYR	2.3
1	C	289	TYR	2.3
2	D	56	GLY	2.3
2	D	54	ASP	2.3
1	C	332	PHE	2.3
2	D	111	ARG	2.2
2	B	246	ILE	2.2
2	D	35	ILE	2.2
2	D	177	ASP	2.2
2	D	47	LYS	2.2
2	D	221	TYR	2.2
2	B	64	ALA	2.1
2	D	38	GLY	2.0
2	B	201	GLU	2.0
2	D	13	GLU	2.0
1	C	263	MET	2.0
2	D	235	LEU	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 ⓘ

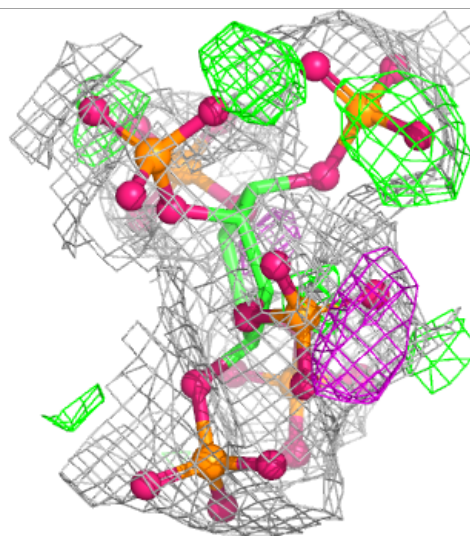
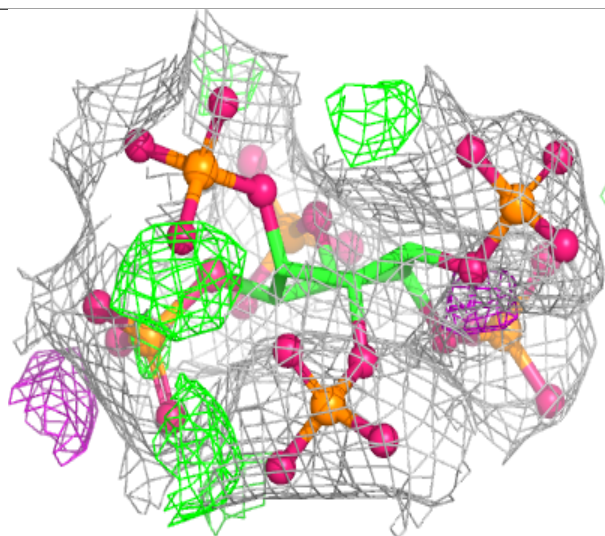
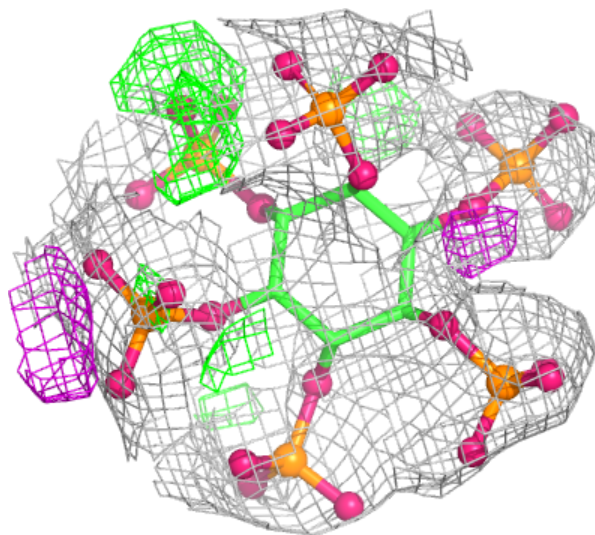
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	IHP	B	402	36/36	0.57	0.16	115,140,144,144	0
3	IHP	D	401	36/36	0.57	0.14	154,166,172,173	0
3	IHP	B	401	36/36	0.60	0.15	166,172,177,177	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.

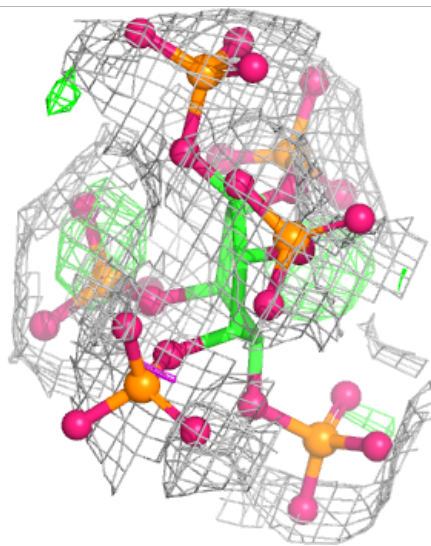
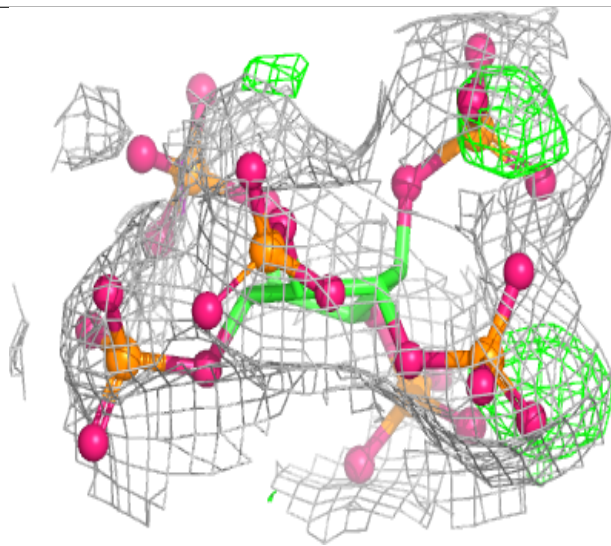
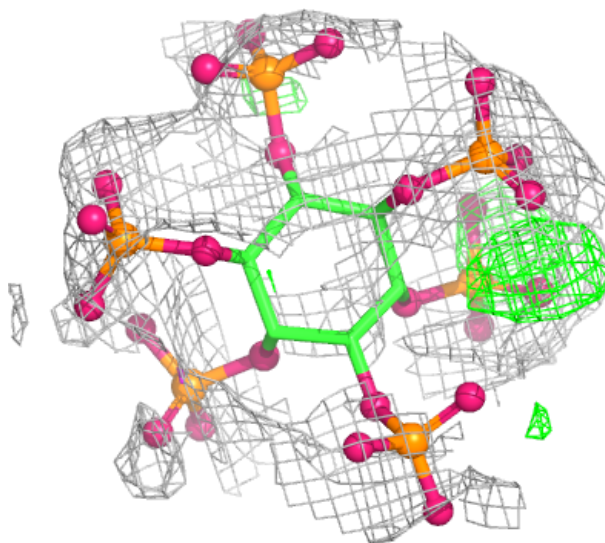
Electron density around IHP B 402:

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



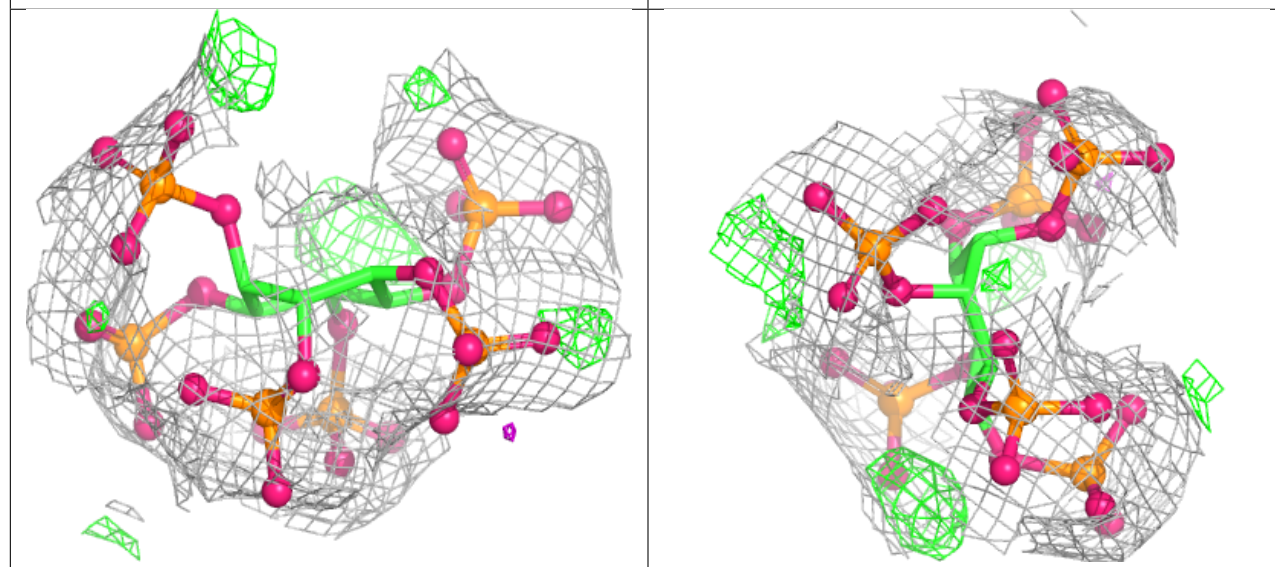
Electron density around IHP D 401:

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



Electron density around IHP B 401:

$2mF_o-DF_c$ (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.