



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 04:10 am BST

PDB ID : 4RXC  
Title : T. Brucei Farnesyl Diphosphate Synthase Complexed with Homorisedronate BPH-6  
Authors : Cao, R.; Liu, Y.-L.; Oldfield, E.  
Deposited on : 2014-12-09  
Resolution : 2.31 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

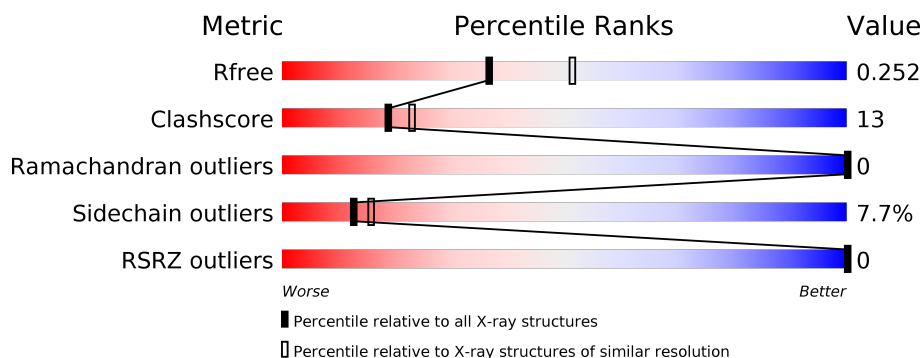
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.31 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	5974 (2.34-2.30)
Clashscore	141614	6604 (2.34-2.30)
Ramachandran outliers	138981	6523 (2.34-2.30)
Sidechain outliers	138945	6523 (2.34-2.30)
RSRZ outliers	127900	5855 (2.34-2.30)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	390	
1	B	390	

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
2	HRX	A	405	-	X	-	-
2	HRX	B	405	-	X	-	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5871 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 Farnesyl pyrophosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	351	Total	C	N	O	S	0	0	0
			2815	1796	464	527	28			
1	B	352	Total	C	N	O	S	0	0	0
			2823	1800	465	530	28			

There are 46 discrepancies between the modelled and reference sequences:

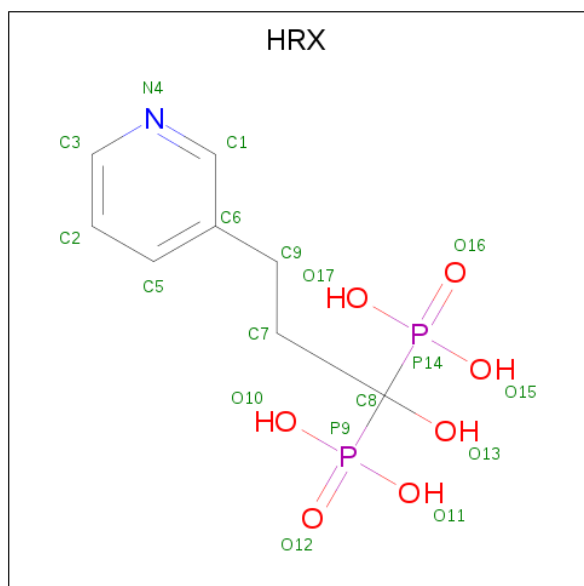
Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	EXPRESSION TAG	UNP Q86C09
A	-21	GLY	-	EXPRESSION TAG	UNP Q86C09
A	-20	SER	-	EXPRESSION TAG	UNP Q86C09
A	-19	SER	-	EXPRESSION TAG	UNP Q86C09
A	-18	HIS	-	EXPRESSION TAG	UNP Q86C09
A	-17	HIS	-	EXPRESSION TAG	UNP Q86C09
A	-16	HIS	-	EXPRESSION TAG	UNP Q86C09
A	-15	HIS	-	EXPRESSION TAG	UNP Q86C09
A	-14	HIS	-	EXPRESSION TAG	UNP Q86C09
A	-13	HIS	-	EXPRESSION TAG	UNP Q86C09
A	-12	SER	-	EXPRESSION TAG	UNP Q86C09
A	-11	SER	-	EXPRESSION TAG	UNP Q86C09
A	-10	GLY	-	EXPRESSION TAG	UNP Q86C09
A	-9	LEU	-	EXPRESSION TAG	UNP Q86C09
A	-8	VAL	-	EXPRESSION TAG	UNP Q86C09
A	-7	PRO	-	EXPRESSION TAG	UNP Q86C09
A	-6	ARG	-	EXPRESSION TAG	UNP Q86C09
A	-5	GLY	-	EXPRESSION TAG	UNP Q86C09
A	-4	SER	-	EXPRESSION TAG	UNP Q86C09
A	-3	HIS	-	EXPRESSION TAG	UNP Q86C09
A	-2	MET	-	EXPRESSION TAG	UNP Q86C09
A	-1	ALA	-	EXPRESSION TAG	UNP Q86C09
A	0	SER	-	EXPRESSION TAG	UNP Q86C09
B	-22	MET	-	EXPRESSION TAG	UNP Q86C09
B	-21	GLY	-	EXPRESSION TAG	UNP Q86C09

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	SER	-	EXPRESSION TAG	UNP Q86C09
B	-19	SER	-	EXPRESSION TAG	UNP Q86C09
B	-18	HIS	-	EXPRESSION TAG	UNP Q86C09
B	-17	HIS	-	EXPRESSION TAG	UNP Q86C09
B	-16	HIS	-	EXPRESSION TAG	UNP Q86C09
B	-15	HIS	-	EXPRESSION TAG	UNP Q86C09
B	-14	HIS	-	EXPRESSION TAG	UNP Q86C09
B	-13	HIS	-	EXPRESSION TAG	UNP Q86C09
B	-12	SER	-	EXPRESSION TAG	UNP Q86C09
B	-11	SER	-	EXPRESSION TAG	UNP Q86C09
B	-10	GLY	-	EXPRESSION TAG	UNP Q86C09
B	-9	LEU	-	EXPRESSION TAG	UNP Q86C09
B	-8	VAL	-	EXPRESSION TAG	UNP Q86C09
B	-7	PRO	-	EXPRESSION TAG	UNP Q86C09
B	-6	ARG	-	EXPRESSION TAG	UNP Q86C09
B	-5	GLY	-	EXPRESSION TAG	UNP Q86C09
B	-4	SER	-	EXPRESSION TAG	UNP Q86C09
B	-3	HIS	-	EXPRESSION TAG	UNP Q86C09
B	-2	MET	-	EXPRESSION TAG	UNP Q86C09
B	-1	ALA	-	EXPRESSION TAG	UNP Q86C09
B	0	SER	-	EXPRESSION TAG	UNP Q86C09

- Molecule 2 is [1-hydroxy-3-(pyridin-3-yl)propane-1,1-diyl]bis(phosphonic acid) (three-letter code: HRX) (formula:  $C_8H_{13}NO_7P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			18	8	1	7	2		
2	A	1	Total	C	N	O	P	0	0
			18	8	1	7	2		
2	B	1	Total	C	N	O	P	0	0
			18	8	1	7	2		
2	B	1	Total	C	N	O	P	0	0
			18	8	1	7	2		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	4	Total	Mg	0	0
			4	4		
3	A	4	Total	Mg	0	0
			4	4		

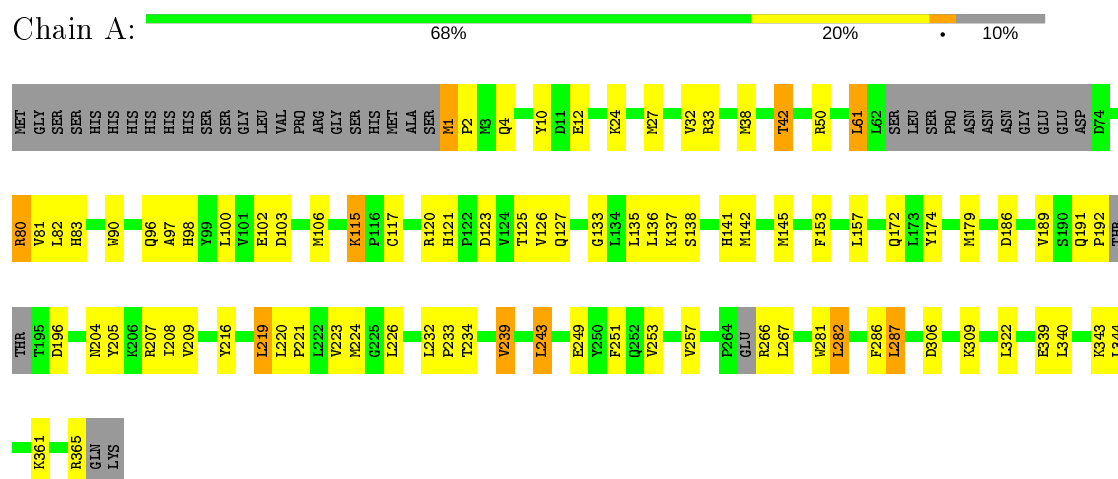
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	84	Total	O	0	0
			84	84		
4	B	69	Total	O	0	0
			69	69		

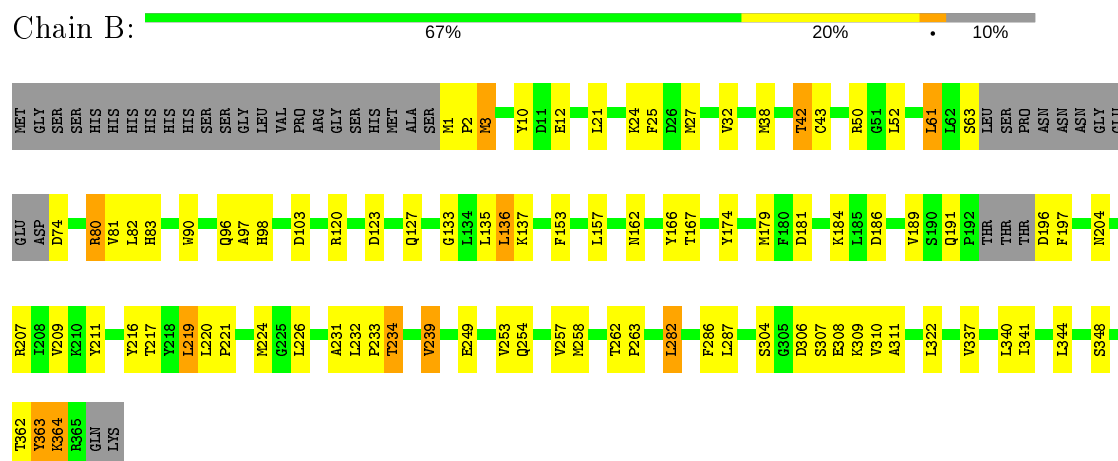
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Farnesyl pyrophosphate synthase



- Molecule 1: Farnesyl pyrophosphate synthase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	133.01Å 120.92Å 63.34Å 90.00° 118.33° 90.00°	Depositor
Resolution (Å)	40.99 – 2.31 40.99 – 2.31	Depositor EDS
% Data completeness (in resolution range)	96.4 (40.99-2.31) 96.4 (40.99-2.31)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.48 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.197 , 0.254 0.201 , 0.252	Depositor DCC
$R_{free}$ test set	1866 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	37.1	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 17.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for k+l,h+l,-l 0.000 for -k+l,-h-l,-l 0.458 for -h-2*l,-k,l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5871	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.96% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: MG, HRX

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	0.21	0/2872	0.32	0/3881
1	B	0.21	0/2881	0.32	0/3894
All	All	0.21	0/5753	0.32	0/7775

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	B	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	362	THR	Peptide
1	B	363	TYR	Peptide
1	B	364	LYS	Peptide

### 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2815	0	2773	84	0
1	B	2823	0	2778	76	0
2	A	36	0	20	4	0
2	B	36	0	19	2	0
3	A	4	0	0	0	0
3	B	4	0	0	1	0
4	A	84	0	0	8	0
4	B	69	0	0	5	0
All	All	5871	0	5590	148	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 13.

All (148) 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:33:ARG:HD3	4:A:501:HOH:O	1.62	1.00
1:A:186:ASP:O	1:B:123:ASP:CB	2.11	0.99
1:A:123:ASP:HB2	1:B:189:VAL:O	1.67	0.95
1:A:103:ASP:HA	1:A:106:MET:HE3	1.51	0.92
1:B:98:HIS:CE1	1:B:137:LYS:HE3	2.05	0.91
1:B:38:MET:O	1:B:42:THR:HB	1.73	0.89
3:B:402:MG:MG	4:B:548:HOH:O	1.18	0.86
1:A:42:THR:HG21	1:A:136:LEU:HD11	1.58	0.85
1:A:103:ASP:HA	1:A:106:MET:CE	2.09	0.82
1:A:186:ASP:O	1:B:123:ASP:HB3	1.79	0.82
1:B:307:SER:O	1:B:310:VAL:HG12	1.79	0.82
1:A:115:LYS:N	1:A:115:LYS:HD2	1.96	0.81
1:A:186:ASP:O	1:B:123:ASP:HB2	1.80	0.79
1:A:123:ASP:CB	1:B:186:ASP:O	2.32	0.78
1:A:123:ASP:HB3	1:B:186:ASP:O	1.85	0.77
1:A:42:THR:HG21	1:A:136:LEU:CD1	2.16	0.75
1:A:205:TYR:O	1:A:208:ILE:HG13	1.88	0.74
1:B:231:ALA:O	1:B:234:THR:HB	1.88	0.73
1:A:33:ARG:NH2	4:A:548:HOH:O	2.20	0.73
1:A:145:MET:HE1	1:B:162:ASN:CB	2.19	0.72
1:A:145:MET:HE1	1:B:162:ASN:HB3	1.72	0.72
1:A:189:VAL:O	1:B:123:ASP:HB2	1.89	0.71
1:A:239:VAL:CG2	1:A:344:LEU:HD13	2.20	0.71
1:A:12:GLU:OE2	1:A:83:HIS:HE1	1.71	0.70
1:A:141:HIS:HB3	1:A:145:MET:CE	2.21	0.70
1:B:196:ASP:O	1:B:197:PHE:HB2	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:239:VAL:CG2	1:B:344:LEU:HD13	2.22	0.69
1:B:1:MET:HB3	1:B:2:PRO:HD3	1.76	0.66
1:A:42:THR:HG23	1:A:97:ALA:HB1	1.77	0.66
1:A:42:THR:CG2	1:A:97:ALA:HB1	2.25	0.66
1:B:103:ASP:OD2	4:B:548:HOH:O	2.15	0.63
1:B:3:MET:HA	1:B:3:MET:CE	2.29	0.63
1:A:208:ILE:HG13	1:A:209:VAL:N	2.14	0.62
1:B:12:GLU:OE2	1:B:83:HIS:HE1	1.83	0.62
1:A:10:TYR:HB2	1:A:90:TRP:CZ2	2.35	0.61
1:A:205:TYR:O	1:A:208:ILE:CG1	2.48	0.60
1:A:239:VAL:HG21	1:A:344:LEU:HD13	1.82	0.59
1:B:1:MET:HB3	1:B:2:PRO:CD	2.32	0.59
1:A:234:THR:HG22	4:A:560:HOH:O	2.03	0.59
1:A:208:ILE:HG13	1:A:209:VAL:H	1.68	0.58
1:B:337:VAL:O	1:B:341:ILE:HG13	2.03	0.58
1:B:3:MET:HE3	1:B:52:LEU:HD21	1.85	0.58
1:B:74:ASP:HB3	4:B:559:HOH:O	2.05	0.57
1:B:181:ASP:HB3	1:B:184:LYS:HG3	1.87	0.57
1:A:125:THR:HG22	1:A:127:GLN:H	1.69	0.57
1:A:196:ASP:C	1:A:196:ASP:OD1	2.43	0.57
1:A:257:VAL:HG22	1:A:282:LEU:HD11	1.85	0.57
1:B:258:MET:O	1:B:262:THR:OG1	2.22	0.57
1:B:239:VAL:HG21	1:B:344:LEU:HD13	1.86	0.56
1:B:10:TYR:HB2	1:B:90:TRP:CZ2	2.41	0.56
1:B:204:ASN:ND2	4:B:566:HOH:O	2.39	0.56
1:B:42:THR:HG21	1:B:136:LEU:HD21	1.87	0.55
1:A:38:MET:O	1:A:42:THR:HB	2.07	0.55
1:A:27:MET:HG2	1:A:32:VAL:HG23	1.89	0.55
1:A:141:HIS:HB3	1:A:145:MET:HE2	1.87	0.54
1:A:232:LEU:N	1:A:233:PRO:CD	2.69	0.54
1:A:234:THR:CG2	4:A:560:HOH:O	2.55	0.54
1:B:157:LEU:HD11	1:B:224:MET:HE3	1.90	0.54
1:A:80:ARG:HG3	1:A:81:VAL:N	2.22	0.54
1:B:254:GLN:HE22	1:B:364:LYS:HA	1.73	0.53
1:B:50:ARG:CZ	1:B:96:GLN:HE22	2.21	0.53
2:A:405:HRX:O12	2:A:405:HRX:O17	2.23	0.53
1:A:61:LEU:HD21	1:A:226:LEU:HD23	1.91	0.53
1:A:306:ASP:HB3	1:A:309:LYS:HG3	1.90	0.53
1:A:157:LEU:HD11	1:A:224:MET:CE	2.39	0.53
2:A:401:HRX:H9	2:A:401:HRX:O17	2.08	0.53
1:A:172:GLN:CD	2:A:401:HRX:H10	2.30	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:ARG:HG3	1:B:81:VAL:N	2.24	0.51
1:B:167:THR:HG23	1:B:211:TYR:HD1	1.75	0.51
1:B:232:LEU:N	1:B:233:PRO:CD	2.73	0.51
1:B:3:MET:HE3	1:B:3:MET:HA	1.93	0.51
1:A:220:LEU:N	1:A:221:PRO:CD	2.73	0.51
1:A:123:ASP:CB	1:B:189:VAL:O	2.51	0.50
1:A:157:LEU:O	1:A:157:LEU:HD12	2.11	0.50
1:B:157:LEU:HD11	1:B:224:MET:CE	2.42	0.49
1:B:174:TYR:CG	1:B:207:ARG:HD3	2.48	0.49
1:B:220:LEU:N	1:B:221:PRO:CD	2.75	0.49
1:B:310:VAL:HG13	1:B:311:ALA:N	2.28	0.49
1:A:117:CYS:O	1:A:121:HIS:HD2	1.97	0.48
1:A:123:ASP:HB2	1:B:186:ASP:O	2.09	0.48
1:A:102:GLU:O	1:A:106:MET:HG3	2.13	0.48
1:A:186:ASP:C	1:B:123:ASP:HB3	2.34	0.48
1:A:189:VAL:HG11	4:A:552:HOH:O	2.13	0.47
1:B:98:HIS:HD2	1:B:133:GLY:O	1.97	0.47
1:A:157:LEU:HD11	1:A:224:MET:HE3	1.96	0.47
1:B:98:HIS:CE1	1:B:137:LYS:CE	2.90	0.47
1:B:249:GLU:O	1:B:253:VAL:HG23	2.13	0.47
1:A:50:ARG:CZ	1:A:96:GLN:HE22	2.28	0.47
1:A:219:LEU:HD22	1:A:223:VAL:HG23	1.97	0.47
1:A:103:ASP:HA	1:A:106:MET:HE2	1.94	0.46
1:B:310:VAL:CG1	1:B:311:ALA:N	2.77	0.46
1:B:3:MET:HE2	1:B:3:MET:HA	1.96	0.46
1:B:220:LEU:HB3	1:B:221:PRO:HD3	1.96	0.46
1:B:61:LEU:HD21	1:B:226:LEU:HD23	1.97	0.46
1:B:174:TYR:CD1	1:B:207:ARG:HD3	2.50	0.46
1:A:204:ASN:ND2	4:A:562:HOH:O	2.49	0.46
1:B:209:VAL:HG11	1:B:249:GLU:HA	1.97	0.46
1:A:192:PRO:C	4:A:581:HOH:O	2.53	0.46
1:A:253:VAL:HG13	1:A:281:TRP:CG	2.51	0.46
1:B:42:THR:HG22	1:B:43:CYS:SG	2.56	0.46
1:A:174:TYR:CG	1:A:207:ARG:HD3	2.51	0.46
1:A:232:LEU:HB3	1:A:233:PRO:HD3	1.97	0.45
1:B:27:MET:HG2	1:B:32:VAL:HG23	1.98	0.45
1:A:96:GLN:O	1:A:100:LEU:HG	2.17	0.45
1:A:1:MET:HB3	1:A:2:PRO:HD3	1.97	0.45
1:A:286:PHE:HB2	1:A:322:LEU:HD11	1.99	0.45
1:A:153:PHE:C	1:A:153:PHE:CD1	2.91	0.45
1:A:189:VAL:O	1:B:123:ASP:OD2	2.34	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:LEU:HA	1:A:61:LEU:HD12	1.83	0.44
1:B:232:LEU:HB3	1:B:233:PRO:HD3	1.98	0.44
1:A:103:ASP:CA	1:A:106:MET:CE	2.90	0.44
1:A:98:HIS:CE1	1:A:137:LYS:NZ	2.86	0.44
1:B:50:ARG:NE	1:B:217:THR:HG23	2.33	0.44
1:A:287:LEU:HD12	1:A:287:LEU:HA	1.78	0.44
1:B:306:ASP:HB3	1:B:309:LYS:HG3	1.99	0.44
1:A:232:LEU:HD23	1:A:232:LEU:C	2.38	0.43
1:B:153:PHE:CD1	1:B:153:PHE:C	2.92	0.43
1:A:102:GLU:C	1:A:106:MET:HE2	2.38	0.43
1:A:243:LEU:HD13	1:A:243:LEU:HA	1.88	0.43
1:A:249:GLU:O	1:A:253:VAL:HG23	2.19	0.43
1:B:282:LEU:HA	1:B:282:LEU:HD12	1.83	0.43
1:A:189:VAL:CG1	4:A:552:HOH:O	2.66	0.42
2:B:405:HRX:H13	4:B:560:HOH:O	2.17	0.42
1:B:42:THR:HG23	1:B:97:ALA:HB1	2.01	0.42
1:B:10:TYR:HB2	1:B:90:TRP:CE2	2.54	0.42
1:B:262:THR:HA	1:B:263:PRO:HD3	1.86	0.42
1:B:363:TYR:O	1:B:363:TYR:CG	2.73	0.42
1:A:106:MET:HB3	1:B:127:GLN:HG3	2.02	0.42
1:B:219:LEU:HA	1:B:219:LEU:HD23	1.83	0.42
1:B:61:LEU:CD2	1:B:226:LEU:HD23	2.50	0.42
1:A:189:VAL:O	1:B:123:ASP:CB	2.65	0.42
1:B:286:PHE:HB2	1:B:322:LEU:HD11	2.02	0.42
1:A:98:HIS:HD2	1:A:133:GLY:O	2.03	0.41
1:A:142:MET:HE3	1:B:166:TYR:HB2	2.02	0.41
1:A:251:PHE:HZ	2:A:405:HRX:C5	2.33	0.41
1:A:174:TYR:CD1	1:A:207:ARG:HD3	2.56	0.41
1:B:98:HIS:ND1	1:B:137:LYS:HE3	2.33	0.41
1:A:239:VAL:CG2	1:A:344:LEU:CD1	2.97	0.41
1:A:1:MET:CB	1:A:2:PRO:HD3	2.51	0.41
1:B:21:LEU:O	1:B:25:PHE:HB2	2.20	0.41
1:B:257:VAL:HG22	1:B:282:LEU:HD11	2.03	0.41
1:A:174:TYR:HB3	1:A:208:ILE:HG22	2.03	0.40
1:B:344:LEU:HD12	1:B:344:LEU:HA	1.88	0.40
1:A:125:THR:HG22	1:A:126:VAL:N	2.36	0.40
1:A:10:TYR:HB2	1:A:90:TRP:CE2	2.56	0.40
1:A:145:MET:HE1	1:B:162:ASN:HB2	1.99	0.40
1:B:308:GLU:H	1:B:308:GLU:HG2	1.73	0.40
2:B:405:HRX:O16	2:B:405:HRX:O11	2.38	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	343/390 (88%)	336 (98%)	7 (2%)	0	100	100
1	B	346/390 (89%)	338 (98%)	8 (2%)	0	100	100
All	All	689/780 (88%)	674 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	306/340 (90%)	280 (92%)	26 (8%)	10	12
1	B	307/340 (90%)	286 (93%)	21 (7%)	16	20
All	All	613/680 (90%)	566 (92%)	47 (8%)	13	16

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	4	GLN
1	A	24	LYS
1	A	42	THR
1	A	61	LEU
1	A	80	ARG
1	A	82	LEU
1	A	115	LYS

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Mol	Chain	Res	Type
1	A	120	ARG
1	A	135	LEU
1	A	138	SER
1	A	179	MET
1	A	191	GLN
1	A	216	TYR
1	A	219	LEU
1	A	239	VAL
1	A	243	LEU
1	A	266	ARG
1	A	267	LEU
1	A	282	LEU
1	A	287	LEU
1	A	339	GLU
1	A	340	LEU
1	A	343	LYS
1	A	361	LYS
1	A	365	ARG
1	B	3	MET
1	B	24	LYS
1	B	42	THR
1	B	61	LEU
1	B	63	SER
1	B	80	ARG
1	B	82	LEU
1	B	120	ARG
1	B	135	LEU
1	B	136	LEU
1	B	179	MET
1	B	191	GLN
1	B	216	TYR
1	B	219	LEU
1	B	234	THR
1	B	239	VAL
1	B	282	LEU
1	B	287	LEU
1	B	304	SER
1	B	340	LEU
1	B	348	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	30	ASN
1	A	83	HIS
1	A	96	GLN
1	A	98	HIS
1	A	121	HIS
1	A	172	GLN
1	A	183	ASN
1	A	204	ASN
1	A	254	GLN
1	A	301	ASN
1	B	30	ASN
1	B	83	HIS
1	B	96	GLN
1	B	98	HIS
1	B	121	HIS
1	B	172	GLN
1	B	183	ASN
1	B	204	ASN
1	B	252	GLN
1	B	254	GLN
1	B	301	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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
2	HRX	A	405	3	18,18,18	2.37	9 (50%)	25,28,28	2.37	11 (44%)
2	HRX	B	405	3	18,18,18	2.59	10 (55%)	25,28,28	2.28	11 (44%)
2	HRX	A	401	3	18,18,18	2.81	10 (55%)	25,28,28	2.30	11 (44%)
2	HRX	B	401	3	18,18,18	2.36	9 (50%)	25,28,28	2.23	9 (36%)

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	HRX	A	405	3	-	19/24/24/24	0/1/1/1
2	HRX	B	405	3	-	13/24/24/24	0/1/1/1
2	HRX	A	401	3	-	2/24/24/24	0/1/1/1
2	HRX	B	401	3	-	5/24/24/24	0/1/1/1

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	HRX	P14-C8	6.19	1.89	1.85
2	B	405	HRX	C9-C6	-4.00	1.40	1.51
2	B	401	HRX	P14-O17	-3.96	1.47	1.54
2	B	401	HRX	P14-O15	-3.95	1.47	1.54
2	B	405	HRX	P14-C8	3.95	1.88	1.85
2	A	405	HRX	C9-C6	-3.94	1.40	1.51
2	A	401	HRX	P14-O17	-3.93	1.47	1.54
2	B	405	HRX	P14-O17	-3.93	1.47	1.54
2	A	405	HRX	P14-O15	-3.91	1.47	1.54
2	A	401	HRX	C9-C6	-3.91	1.40	1.51
2	B	405	HRX	P14-O15	-3.90	1.47	1.54
2	B	401	HRX	C9-C6	-3.90	1.40	1.51
2	B	401	HRX	P9-O10	-3.88	1.47	1.54
2	A	401	HRX	P14-O15	-3.86	1.47	1.54
2	A	405	HRX	P14-O17	-3.85	1.47	1.54
2	A	401	HRX	P9-O10	-3.80	1.47	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	405	HRX	P9-O10	-3.74	1.47	1.54
2	B	405	HRX	P9-O10	-3.66	1.48	1.54
2	A	405	HRX	P9-O11	-3.51	1.48	1.54
2	B	405	HRX	P9-O11	-3.47	1.48	1.54
2	B	401	HRX	P9-O11	-3.47	1.48	1.54
2	A	401	HRX	P9-O11	-3.42	1.48	1.54
2	B	405	HRX	C1-N4	2.93	1.40	1.34
2	A	401	HRX	C1-N4	2.92	1.40	1.34
2	A	405	HRX	C1-N4	2.84	1.40	1.34
2	B	401	HRX	C1-N4	2.79	1.40	1.34
2	B	405	HRX	P9-C8	-2.78	1.83	1.85
2	A	405	HRX	P14-C8	2.57	1.87	1.85
2	A	401	HRX	P9-C8	-2.35	1.83	1.85
2	A	401	HRX	C3-N4	2.26	1.40	1.33
2	B	401	HRX	P14-C8	2.25	1.86	1.85
2	A	405	HRX	C3-N4	2.25	1.40	1.33
2	B	405	HRX	C3-N4	2.25	1.40	1.33
2	B	401	HRX	C3-N4	2.19	1.40	1.33
2	B	405	HRX	P9-O12	-2.14	1.47	1.50
2	B	401	HRX	P9-O12	-2.10	1.47	1.50
2	A	401	HRX	P14-O16	-2.07	1.47	1.50
2	A	405	HRX	P9-O12	-2.01	1.47	1.50

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	405	HRX	O16-P14-C8	-5.00	96.77	109.86
2	B	405	HRX	O16-P14-C8	-4.78	97.35	109.86
2	A	401	HRX	O16-P14-C8	-4.68	97.62	109.86
2	B	401	HRX	O16-P14-C8	-4.65	97.69	109.86
2	A	405	HRX	O15-P14-C8	-4.39	96.33	106.17
2	B	405	HRX	O15-P14-C8	-4.22	96.71	106.17
2	B	401	HRX	O17-P14-C8	-4.19	96.78	106.17
2	B	401	HRX	O15-P14-C8	-4.16	96.83	106.17
2	A	401	HRX	O17-P14-C8	-4.12	96.93	106.17
2	A	405	HRX	O17-P14-C8	-3.85	97.54	106.17
2	A	405	HRX	C7-C9-C6	-3.83	102.82	112.85
2	B	405	HRX	O15-P14-O17	3.82	118.77	107.99
2	A	401	HRX	O15-P14-C8	-3.81	97.61	106.17
2	A	401	HRX	O15-P14-O17	3.80	118.71	107.99
2	B	405	HRX	O17-P14-C8	-3.75	97.75	106.17
2	B	401	HRX	O15-P14-O17	3.70	118.44	107.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	405	HRX	O15-P14-O17	3.66	118.33	107.99
2	A	401	HRX	C7-C9-C6	-3.62	103.36	112.85
2	B	405	HRX	C7-C9-C6	-3.14	104.63	112.85
2	A	405	HRX	O15-P14-O16	3.02	119.83	113.06
2	B	401	HRX	O15-P14-O16	2.98	119.74	113.06
2	A	401	HRX	O11-P9-C8	-2.78	99.94	106.17
2	B	405	HRX	O17-P14-O16	2.71	119.14	113.06
2	A	405	HRX	O10-P9-O11	2.65	115.47	107.99
2	B	405	HRX	O10-P9-O11	2.58	115.29	107.99
2	A	405	HRX	P9-C8-P14	-2.52	108.30	112.81
2	A	405	HRX	O12-P9-C8	-2.47	103.39	109.86
2	A	401	HRX	O17-P14-O16	2.45	118.55	113.06
2	A	401	HRX	O10-P9-O11	2.40	114.77	107.99
2	A	401	HRX	O12-P9-C8	-2.32	103.78	109.86
2	B	401	HRX	C9-C6-C1	-2.30	118.13	121.83
2	B	401	HRX	O10-P9-O11	2.26	114.36	107.99
2	A	405	HRX	O11-P9-C8	-2.21	101.22	106.17
2	A	401	HRX	O15-P14-O16	2.14	117.86	113.06
2	A	401	HRX	P9-C8-P14	-2.11	109.03	112.81
2	B	405	HRX	P9-C8-P14	-2.07	109.11	112.81
2	B	405	HRX	O12-P9-C8	-2.06	104.46	109.86
2	B	401	HRX	C7-C9-C6	-2.06	107.47	112.85
2	B	401	HRX	O11-P9-C8	-2.03	101.61	106.17
2	B	405	HRX	O11-P9-C8	-2.03	101.61	106.17
2	B	405	HRX	O10-P9-C8	-2.02	101.63	106.17
2	A	405	HRX	O17-P14-O16	2.02	117.59	113.06

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	405	HRX	P9-C8-P14-O17
2	A	405	HRX	O13-C8-P14-O17
2	A	405	HRX	C7-C8-P14-O17
2	A	405	HRX	P9-C8-P14-O16
2	A	405	HRX	O13-C8-P14-O16
2	A	405	HRX	C7-C8-P14-O16
2	A	405	HRX	P9-C8-P14-O15
2	A	405	HRX	O13-C8-P14-O15
2	A	405	HRX	C7-C8-P14-O15
2	A	405	HRX	O13-C8-P9-O12
2	B	405	HRX	P9-C8-P14-O17

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Mol	Chain	Res	Type	Atoms
2	B	405	HRX	O13-C8-P14-O17
2	B	405	HRX	C7-C8-P14-O17
2	B	405	HRX	P9-C8-P14-O16
2	B	405	HRX	O13-C8-P14-O16
2	B	405	HRX	C7-C8-P14-O16
2	B	405	HRX	P9-C8-P14-O15
2	B	405	HRX	O13-C8-P14-O15
2	B	405	HRX	C7-C8-P14-O15
2	B	405	HRX	C9-C7-C8-P14
2	B	405	HRX	C9-C7-C8-P9
2	B	405	HRX	C9-C7-C8-O13
2	A	401	HRX	C8-C7-C9-C6
2	B	401	HRX	C9-C7-C8-P14
2	B	401	HRX	C9-C7-C8-P9
2	B	401	HRX	C9-C7-C8-O13
2	A	405	HRX	C9-C7-C8-P14
2	A	405	HRX	C9-C7-C8-P9
2	A	405	HRX	C9-C7-C8-O13
2	A	405	HRX	P14-C8-P9-O12
2	A	405	HRX	C7-C8-P9-O12
2	B	401	HRX	C5-C6-C9-C7
2	A	405	HRX	C5-C6-C9-C7
2	B	405	HRX	C5-C6-C9-C7
2	A	405	HRX	P14-C8-P9-O11
2	A	405	HRX	C7-C8-P9-O10
2	A	401	HRX	C7-C8-P14-O15
2	B	401	HRX	C1-C6-C9-C7
2	A	405	HRX	O13-C8-P9-O10

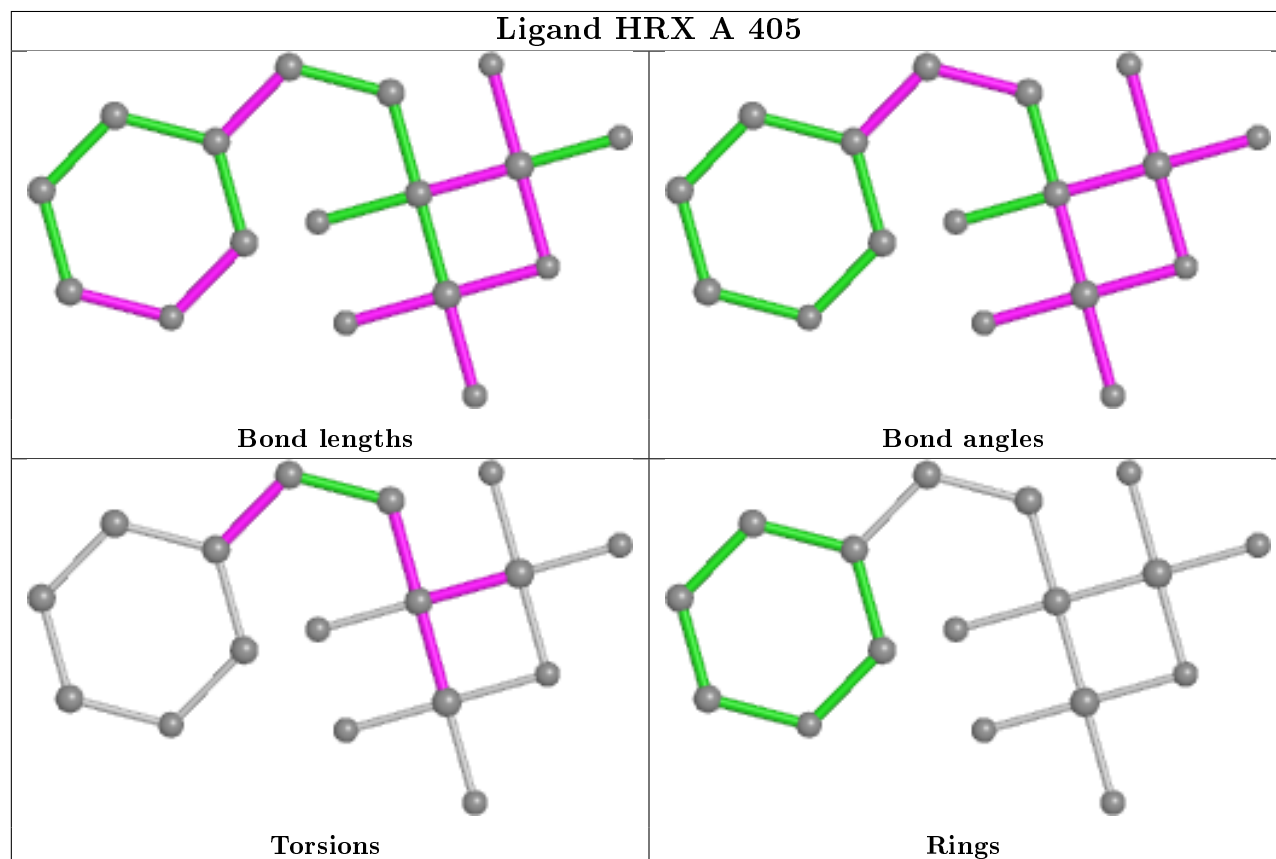
There are no ring outliers.

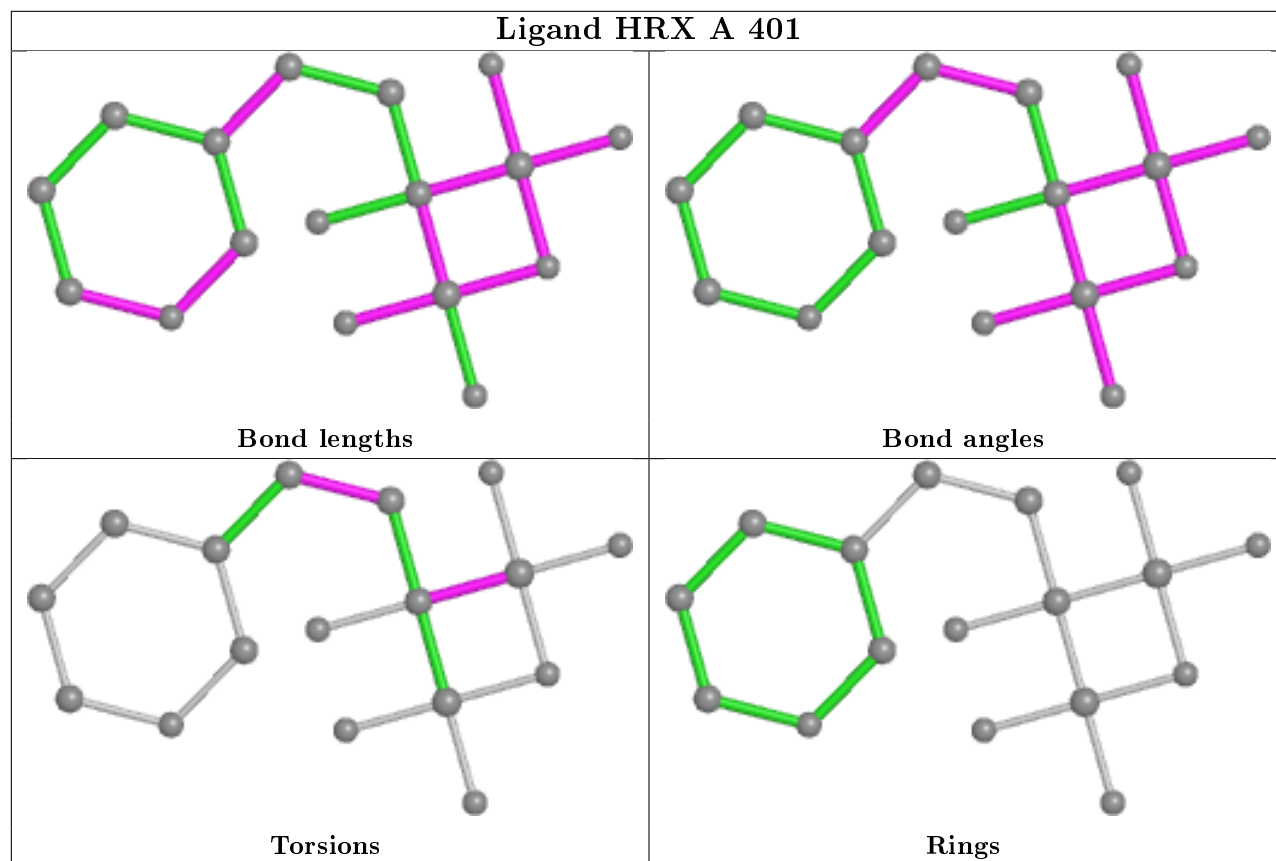
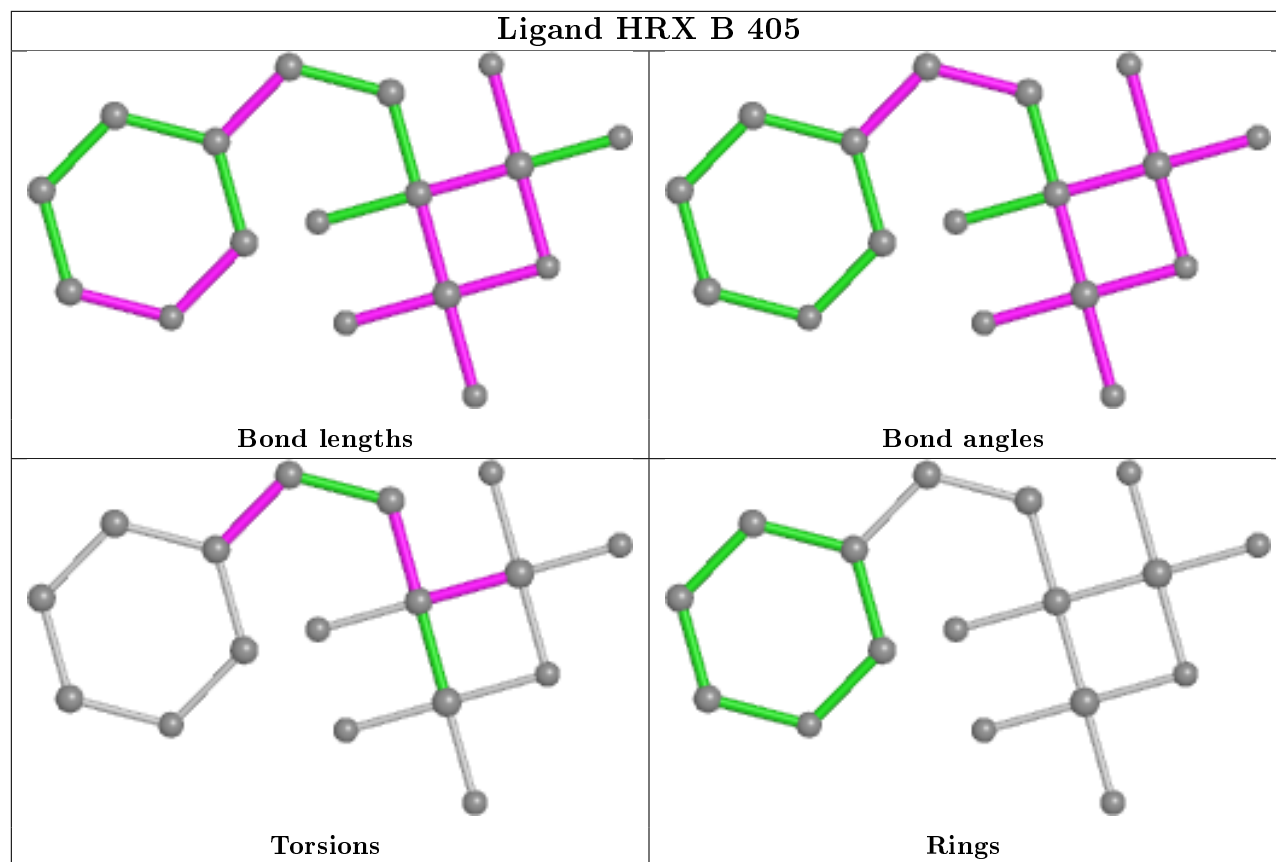
3 monomers are involved in 6 short contacts:

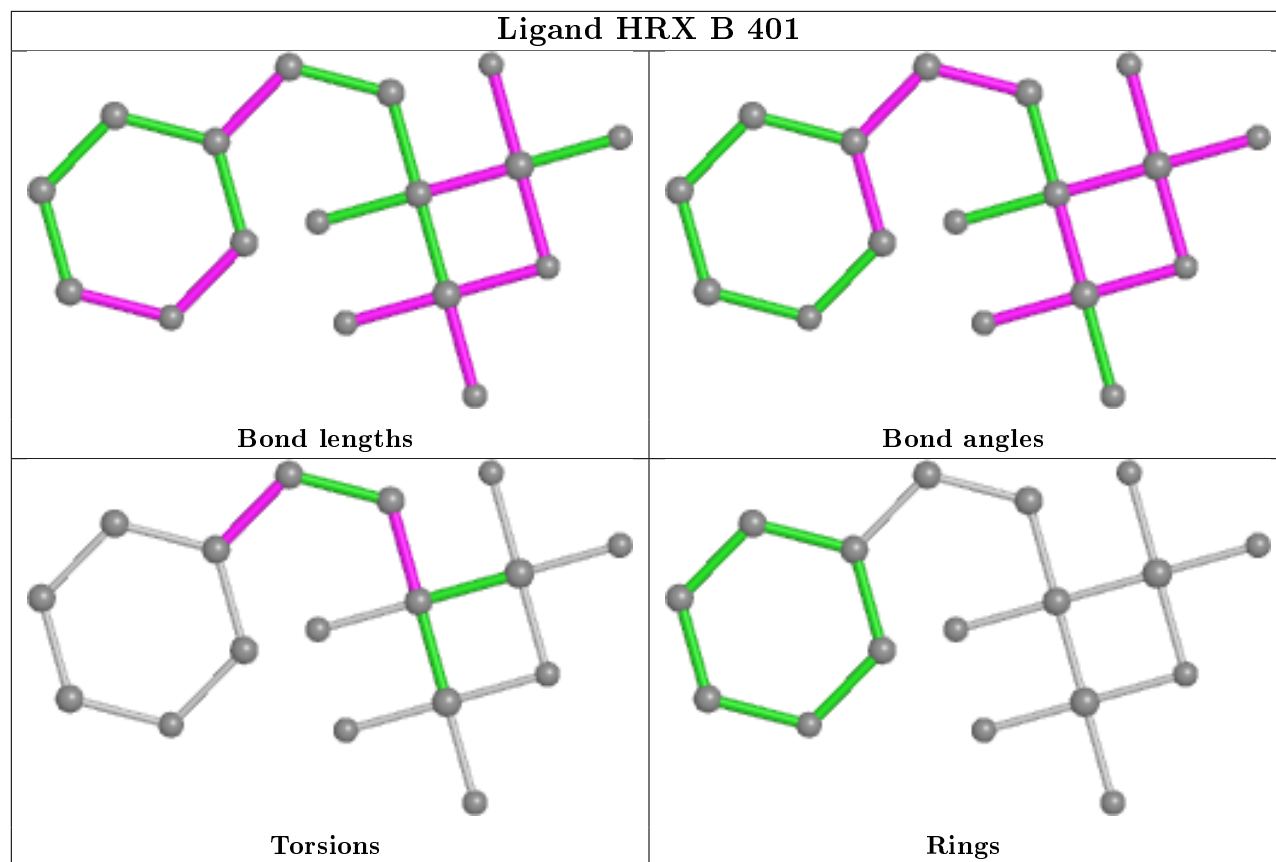
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	405	HRX	2	0
2	B	405	HRX	2	0
2	A	401	HRX	2	0

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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	351/390 (90%)	-0.59	0 100 100	26, 39, 66, 91	0
1	B	352/390 (90%)	-0.60	0 100 100	25, 39, 67, 102	0
All	All	703/780 (90%)	-0.59	0 100 100	25, 39, 67, 102	0

There are no RSRZ outliers to report.

### 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 carbohydrates 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	MG	A	404	1/1	0.81	0.20	49,49,49,49	0
3	MG	B	403	1/1	0.82	0.23	39,39,39,39	0
3	MG	A	406	1/1	0.83	0.24	46,46,46,46	0
3	MG	B	406	1/1	0.85	0.22	47,47,47,47	0
3	MG	A	403	1/1	0.95	0.25	39,39,39,39	0
3	MG	A	402	1/1	0.95	0.18	39,39,39,39	0
2	HRX	A	401	18/18	0.96	0.11	24,38,45,47	0

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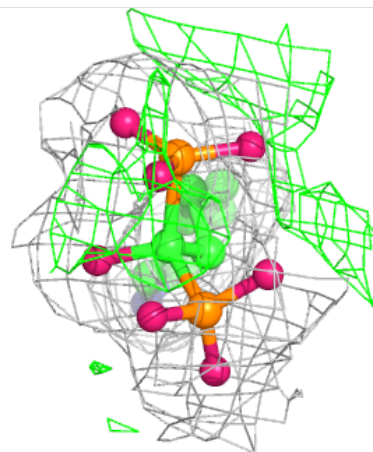
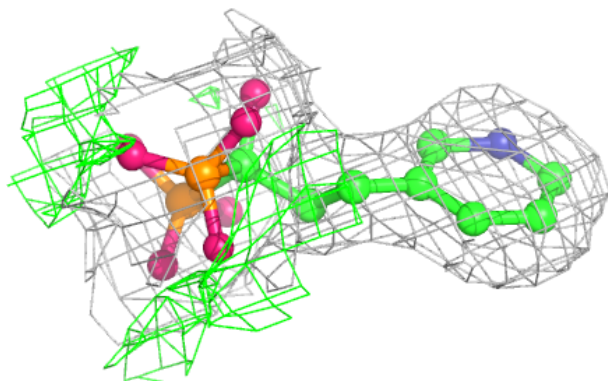
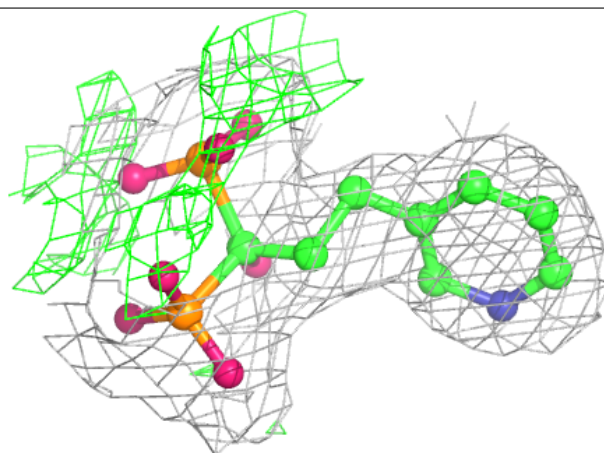
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	HRX	A	405	18/18	0.96	0.13	35,59,101,104	0
3	MG	B	404	1/1	0.96	0.27	53,53,53,53	0
2	HRX	B	405	18/18	0.97	0.10	30,49,91,96	0
3	MG	B	402	1/1	0.97	0.16	55,55,55,55	0
2	HRX	B	401	18/18	0.98	0.11	26,41,50,53	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.

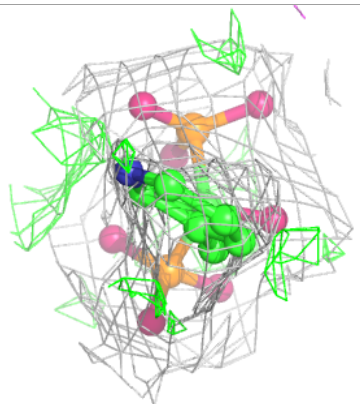
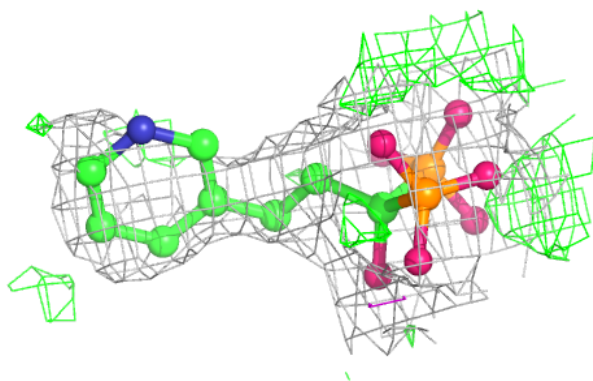
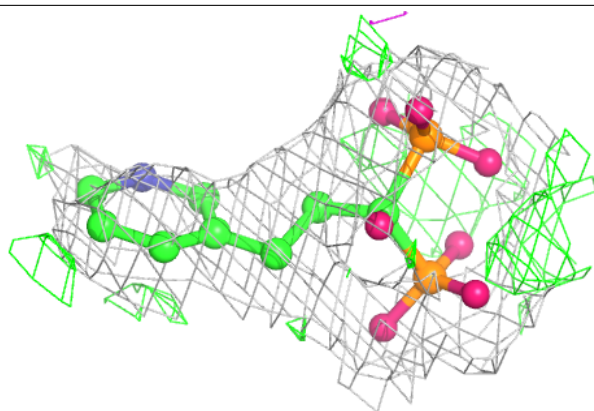
**Electron density around HRX A 401:**

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)

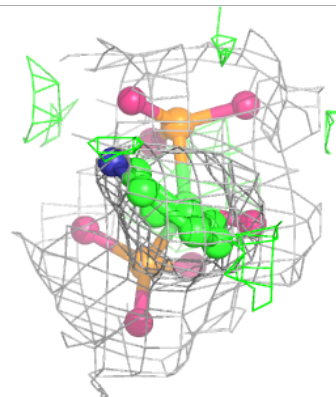
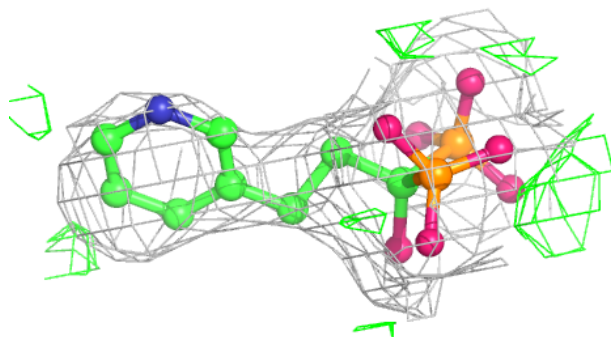
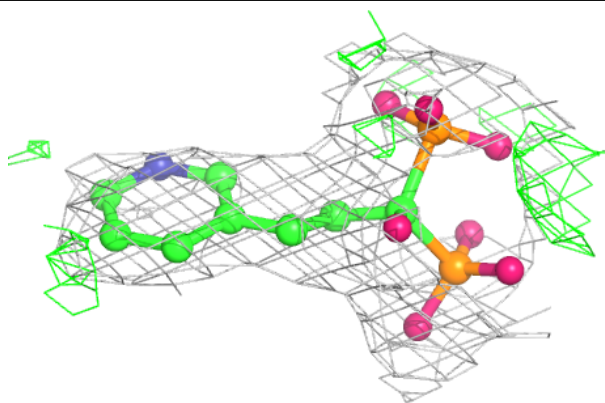


**Electron density around HRX A 405:**

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

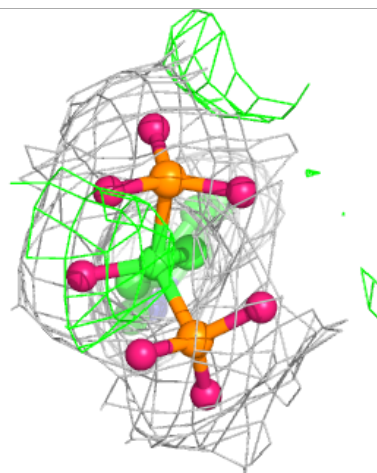
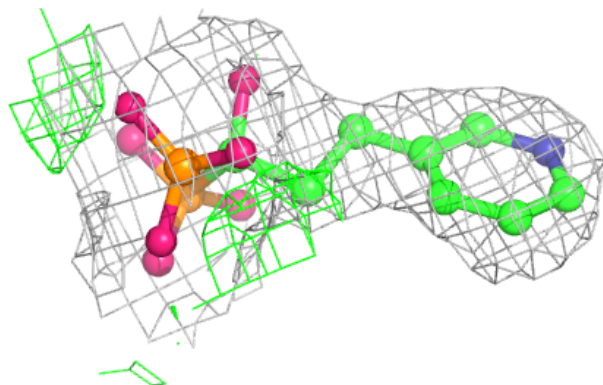
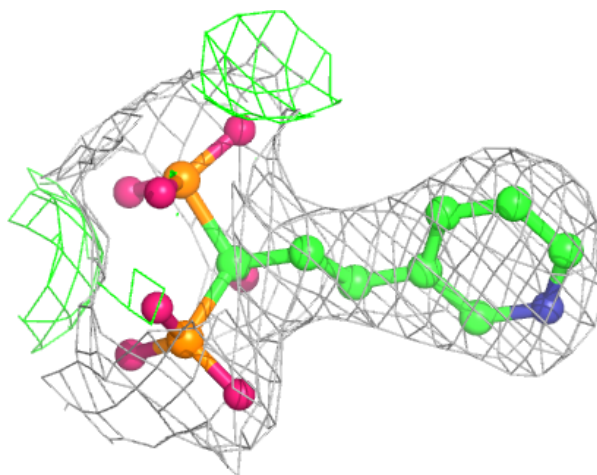
**Electron density around HRX B 405:**

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



**Electron density around HRX 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.