

Full wwPDB X-ray Structure Validation Report (i)

Dec 10, 2022 – 06:28 PM EST

PDB ID : 1A2G

Title : PROBING THE STRENGTH AND CHARACTER OF AN ASP-HIS-X HY-

DROGEN BOND BY INTRODUCING BURIED CHARGES

Authors: Cao, Y.; Goodin, D.B.; Mcree, D.E.

Deposited on : 1998-01-02

Resolution : 2.10 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 (i) 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 (1)) 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 : 2.31.2 buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

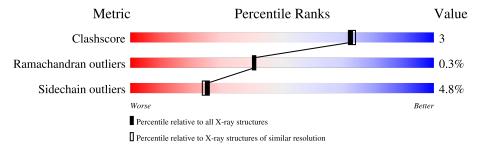
Validation Pipeline (wwPDB-VP) : 2.31.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$

The reported resolution of this entry is 2.10 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)

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 <=5%

Mol	Chain	Length	Quality of chain		
1	A	291	84%	13%	•



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3016 atoms, of which 514 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.

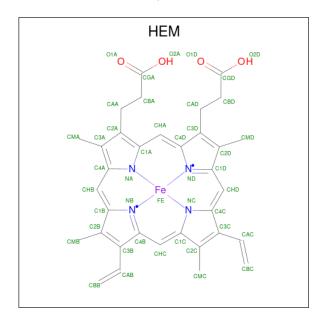
• Molecule 1 is a protein called CYTOCHROME C PEROXIDASE.

Mol	Chain	Residues		Atoms					ZeroOcc	AltConf	Trace
1	A	291	Total	C	H	N	0	S	0	0	0
			2866	1502	514	394	451	5			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	53	ILE	THR	engineered mutation	UNP P00431
A	152	GLY	ASP	engineered mutation	UNP P00431
A	231	HIS	MET	engineered mutation	UNP P00431

• Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	Λ	1	Total	С	Fe	N	О	0	0
2	Α	1	43	34	1	4	4	0	0



• Molecule 3 is water.

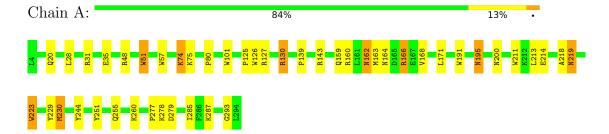
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	107	Total O 107 107	0	0



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. 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. 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: CYTOCHROME C PEROXIDASE





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	104.60Å 74.00Å 45.10Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	5.00 - 2.10	Depositor
resolution (A)	42.71 - 2.00	EDS
% Data completeness	(Not available) $(5.00-2.10)$	Depositor
(in resolution range)	84.4 (42.71-2.00)	EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) > 1$	0.15 (at 2.00Å)	Xtriage
Refinement program	XTALVIEW, X-PLOR 3.1	Depositor
P. P.	0.184 , (Not available)	Depositor
R, R_{free}	0.375 , (Not available)	DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	19.7	Xtriage
Anisotropy	0.569	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.29 , 44.1	EDS
L-test for twinning ²	$ < L > = 0.48, < L^2> = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.80	EDS
Total number of atoms	3016	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.26% of the height of the origin peak. No significant pseudotranslation is detected.

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



¹Intensities estimated from amplitudes.

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: HEM

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).

Mal	Chain	Bond	lengths	Bond angles		
IVIOI	Chain	RMSZ	lengths $\# Z > 5$	RMSZ	# Z > 5	
1	A	0.85	0/2419	1.55	40/3275 (1.2%)	

There are no bond length outliers.

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	31	ARG	NE-CZ-NH2	-11.43	114.58	120.30
1	A	127	ARG	NE-CZ-NH2	-11.19	114.71	120.30
1	A	127	ARG	NE-CZ-NH1	9.34	124.97	120.30
1	A	223	TRP	CD1-CG-CD2	8.73	113.28	106.30
1	A	163	MET	CG-SD-CE	-8.72	86.25	100.20
1	A	51	TRP	CD1-CG-CD2	8.04	112.73	106.30
1	A	101	TRP	CD1-CG-CD2	7.80	112.54	106.30
1	A	143	ARG	NE-CZ-NH1	7.73	124.16	120.30
1	A	223	TRP	CE2-CD2-CG	-7.47	101.33	107.30
1	A	143	ARG	NE-CZ-NH2	-7.17	116.71	120.30
1	A	229	TYR	CB-CG-CD2	-7.16	116.70	121.00
1	A	211	TRP	CD1-CG-CD2	7.16	112.03	106.30
1	A	51	TRP	CE2-CD2-CG	-7.07	101.64	107.30
1	A	57	TRP	CD1-CG-CD2	7.03	111.92	106.30
1	A	31	ARG	NE-CZ-NH1	7.01	123.81	120.30
1	A	126	TRP	CD1-CG-CD2	6.94	111.86	106.30
1	A	101	TRP	CE2-CD2-CG	-6.91	101.77	107.30
1	A	191	TRP	CB-CG-CD1	-6.86	118.08	127.00
1	A	191	TRP	CE2-CD2-CG	-6.70	101.94	107.30
1	A	160	ARG	NE-CZ-NH2	-6.64	116.98	120.30
1	A	57	TRP	CE2-CD2-CG	-6.60	102.02	107.30
1	A	211	TRP	CE2-CD2-CG	-6.49	102.10	107.30
1	A	48	ARG	NE-CZ-NH1	6.45	123.53	120.30
1	A	244	TYR	CB-CG-CD2	-6.35	117.19	121.00

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Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	162	ASN	CA-C-N	-6.17	103.62	117.20
1	A	126	TRP	CE2-CD2-CG	-6.11	102.41	107.30
1	A	51	TRP	CB-CG-CD1	-5.89	119.34	127.00
1	A	191	TRP	CD1-CG-CD2	5.86	110.99	106.30
1	A	130	ARG	CG-CD-NE	-5.73	99.76	111.80
1	A	223	TRP	CG-CD1-NE1	-5.66	104.44	110.10
1	A	51	TRP	CG-CD1-NE1	-5.55	104.55	110.10
1	A	230	MET	CG-SD-CE	-5.50	91.40	100.20
1	A	166	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	A	101	TRP	CG-CD1-NE1	-5.29	104.81	110.10
1	A	277	PRO	CA-C-N	5.28	128.82	117.20
1	A	278	LYS	CA-CB-CG	-5.20	101.95	113.40
1	A	251	TYR	CB-CG-CD2	-5.19	117.89	121.00
1	A	191	TRP	CG-CD2-CE3	5.08	138.47	133.90
1	A	162	ASN	CA-C-O	5.08	130.76	120.10
1	A	57	TRP	CG-CD1-NE1	-5.01	105.09	110.10

There are no chirality outliers.

There are no planarity outliers.

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	2352	514	2226	16	2
2	A	43	0	30	0	0
3	A	107	0	0	3	2
All	All	2502	514	2256	16	2

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 3.

All (16) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	1100111 1		$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:A:200:ASN:H	1:A:255:GLN:HE21	1.43	0.67

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COHABABACA		DIEUIUU	DUIUE
0 0 1000100000			

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${f distance}({ m \AA})$	overlap (Å)
1:A:74:LYS:HD2	1:A:74:LYS:H	1.65	0.60
1:A:20:GLN:HE22	1:A:287:LYS:H	1.51	0.58
1:A:125:PRO:HG3	1:A:285:ILE:HD11	1.85	0.58
1:A:130:ARG:NE	3:A:326:HOH:O	2.48	0.46
1:A:125:PRO:HG3	1:A:285:ILE:CD1	2.47	0.44
1:A:75:LYS:HE3	1:A:139:PRO:O	2.18	0.44
1:A:195:ASN:HD22	1:A:195:ASN:H	1.66	0.44
1:A:20:GLN:HE22	1:A:287:LYS:N	2.12	0.43
1:A:164:ASN:O	1:A:168:VAL:HG23	2.19	0.42
1:A:130:ARG:NH2	3:A:326:HOH:O	2.51	0.42
1:A:213:LEU:HD13	1:A:223:TRP:CE2	2.54	0.42
1:A:74:LYS:H	1:A:74:LYS:CD	2.33	0.41
1:A:195:ASN:HB3	3:A:308:HOH:O	2.21	0.41
1:A:218:ALA:O	1:A:219:ASN:HB2	2.21	0.41
1:A:230:MET:SD	1:A:230:MET:C	3.00	0.40

All (2) 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	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:A:214:GLU:OE2	3:A:374:HOH:O[4_4610]	1.92	0.28
1:A:293:GLY:O	3:A:330:HOH:O[3_549]	2.19	0.01

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	289/291 (99%)	276 (96%)	12 (4%)	1 (0%)	41 41	

All (1) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	A	162	ASN

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	Analysed Rotameric Outliers		Percentiles	
1	A	249/249 (100%)	237 (95%)	12 (5%)	25	24

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	35	GLU
1	A	51	TRP
1	A	74	LYS
1	A	80	PRO
1	A	159	GLN
1	A	166	ARG
1	A	171	LEU
1	A	195	ASN
1	A	219	ASN
1	A	260	LYS
1	A	279	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	GLN
1	A	24	ASN
1	A	87	ASN
1	A	195	ASN
1	A	208	ASN
1	A	220	ASN
1	A	231	HIS
1	A	255	GLN
1	A	292	GLN



5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand is 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	B	ond leng	gths	В	ond ang	les
MIOI	OI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	,	HEM	A	1	3,1	41,50,50	1.88	17 (41%)	45,82,82	1.79	7 (15%)

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	HEM	A	1	3,1	-	2/12/54/54	-

All (17) bond length outliers are listed below:

Mo	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(\AA)$	$\operatorname{Ideal}(ext{\AA})$
2	A	1	HEM	C1B-NB	-4.03	1.33	1.40
2	A	1	HEM	C4D-C3D	-3.47	1.39	1.45
2	A	1	HEM	C4D-ND	-3.35	1.34	1.40

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Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(A)
2	A	1	HEM	C3C-CAC	-3.33	1.41	1.47
2	A	1	HEM	C4B-NB	-3.05	1.32	1.38
2	A	1	HEM	CBB-CAB	2.84	1.44	1.30
2	A	1	HEM	C1D-C2D	-2.67	1.39	1.44
2	A	1	HEM	C1D-ND	-2.66	1.33	1.38
2	A	1	HEM	O2A-CGA	-2.47	1.22	1.30
2	A	1	HEM	CAB-C3B	-2.29	1.41	1.47
2	A	1	HEM	C1B-C2B	-2.24	1.40	1.44
2	A	1	HEM	CHA-C4D	2.21	1.40	1.35
2	A	1	HEM	CHB-C1B	2.18	1.40	1.35
2	A	1	HEM	C3B-C4B	-2.17	1.40	1.44
2	A	1	HEM	O2D-CGD	-2.07	1.23	1.30
2	A	1	HEM	FE-NB	2.03	2.06	1.96
2	A	1	HEM	FE-ND	2.01	2.06	1.96

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	A	1	HEM	C4B-CHC-C1C	6.50	131.14	122.56
2	A	1	HEM	C4C-CHD-C1D	5.32	129.58	122.56
2	A	1	HEM	CMC-C2C-C3C	2.97	130.24	124.68
2	A	1	HEM	C3B-C2B-C1B	-2.76	104.44	106.49
2	A	1	HEM	C4B-C3B-C2B	-2.08	105.47	107.11
2	A	1	HEM	CMA-C3A-C4A	-2.05	125.31	128.46
2	A	1	HEM	CMA-C3A-C2A	2.04	128.78	124.94

There are no chirality outliers.

All (2) torsion outliers are listed below:

	Mol	Chain	Res	Type	Atoms
	2	A	1	HEM	CAA-CBA-CGA-O1A
ĺ	2	A	1	HEM	CAA-CBA-CGA-O2A

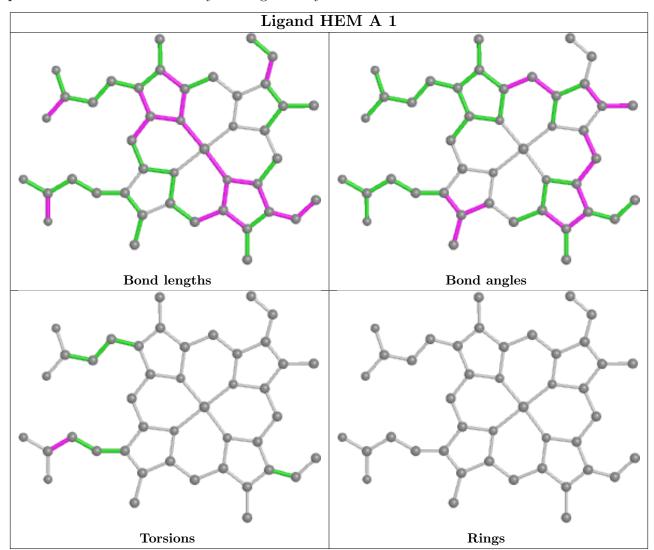
There are no ring outliers.

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 then 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)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

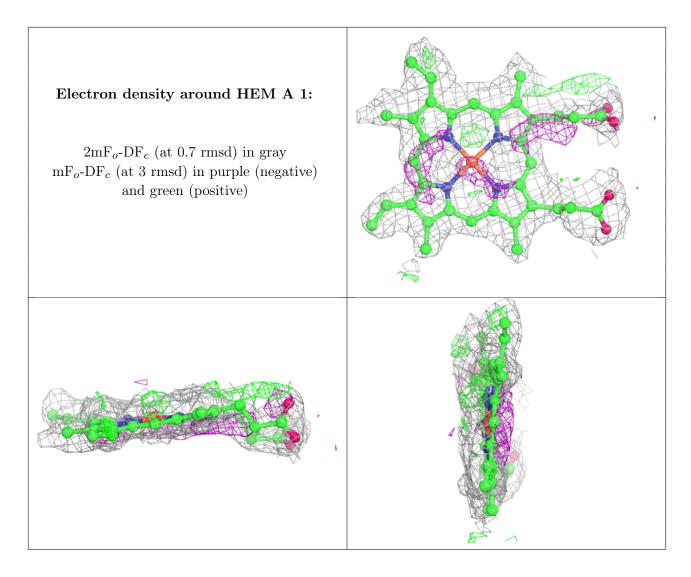
Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

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)

Unable to reproduce the depositors R factor - this section is therefore empty.

