Full wwPDB X-ray Structure Validation Report (i)

Aug 27, 2026 – 07:01 pm BST

PDB ID	:	6Y7D / pdb_00006y7d
Title	:	Pseudomonas stutzeri nitrous oxide reductase mutant, H433A
Authors	:	Zhang, L.; Kroneck, P.M.H.; Einsle, O.
Deposited on	:	2020-02-28
Resolution	:	1.60 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 (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 (i)) were used in the production of this report:

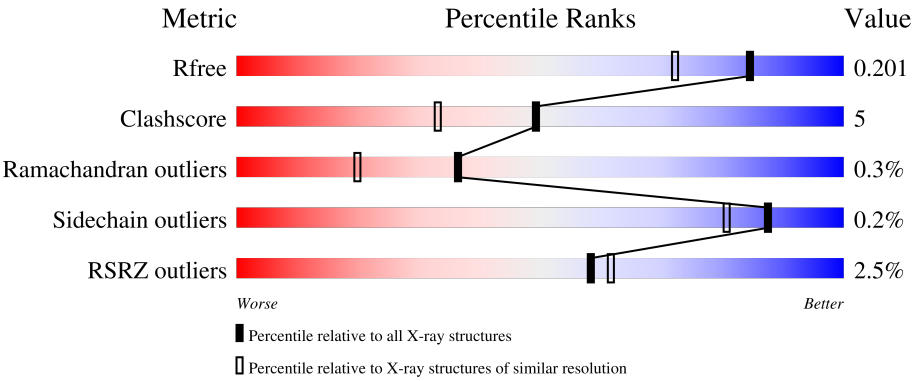
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

1 Overall quality at a glance (i)

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

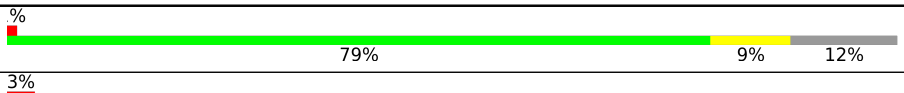
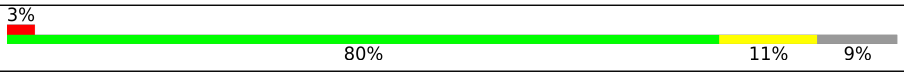
The reported resolution of this entry is 1.60 Å.

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	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

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 <=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	646	
1	B	646	

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 10347 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.

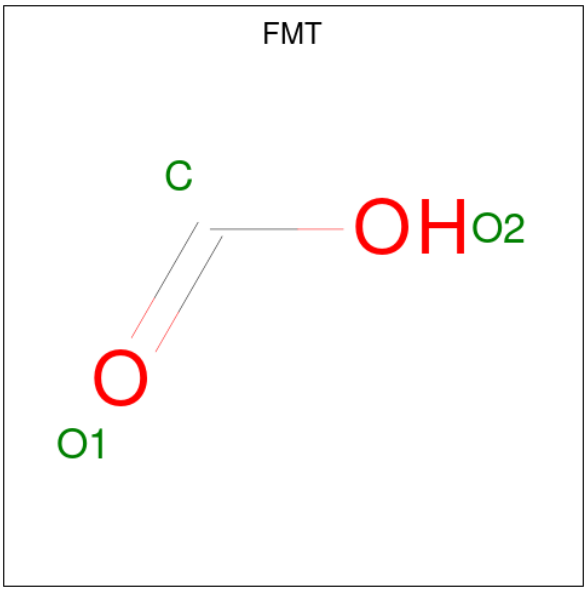
- Molecule 1 is a protein called Nitrous-oxide reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	570	Total	C	N	O	S	0	7	0
			4510	2857	771	852	30			
1	B	588	Total	C	N	O	S	0	2	0
			4634	2933	796	872	33			

There are 18 discrepancies between the modelled and reference sequences:

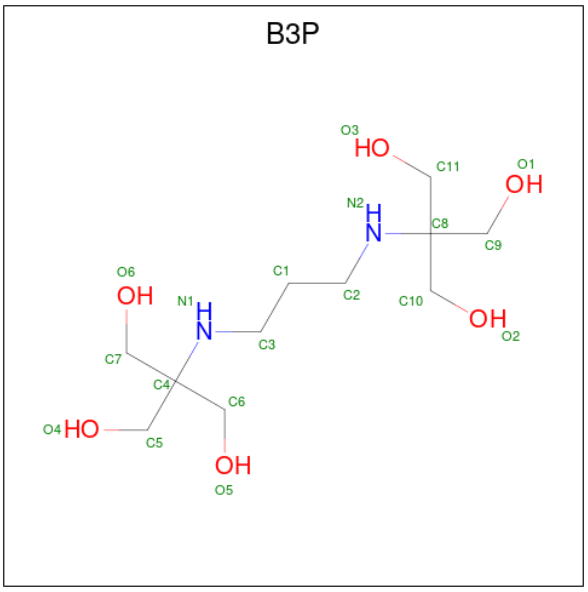
Chain	Residue	Modelled	Actual	Comment	Reference
A	433	ALA	HIS	engineered mutation	UNP P19573
A	639	TRP	-	expression tag	UNP P19573
A	640	SER	-	expression tag	UNP P19573
A	641	HIS	-	expression tag	UNP P19573
A	642	PRO	-	expression tag	UNP P19573
A	643	GLN	-	expression tag	UNP P19573
A	644	PHE	-	expression tag	UNP P19573
A	645	GLU	-	expression tag	UNP P19573
A	646	LYS	-	expression tag	UNP P19573
B	433	ALA	HIS	engineered mutation	UNP P19573
B	639	TRP	-	expression tag	UNP P19573
B	640	SER	-	expression tag	UNP P19573
B	641	HIS	-	expression tag	UNP P19573
B	642	PRO	-	expression tag	UNP P19573
B	643	GLN	-	expression tag	UNP P19573
B	644	PHE	-	expression tag	UNP P19573
B	645	GLU	-	expression tag	UNP P19573
B	646	LYS	-	expression tag	UNP P19573

- Molecule 2 is FORMIC ACID (CCD ID: FMT) (formula: CH_2O_2).



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

- Molecule 3 is 2-[3-(2-HYDROXY-1,1-DIHYDROXYMETHYL-ETHYLAMINO)-PROPYLAMINO]-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: B3P) (formula: $C_{11}H_{26}N_2O_6$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 19 11 2 6	0	0
3	B	1	Total C N O 19 11 2 6	0	0

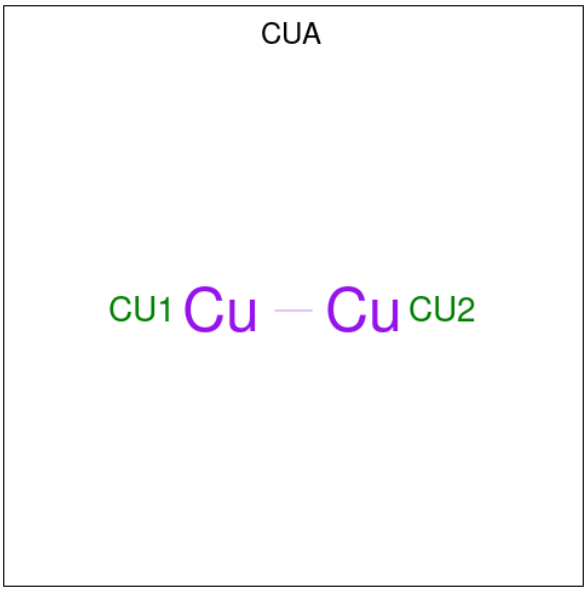
- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Na 1 1	0	0
4	B	1	Total Na 1 1	0	0

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Cl 1 1	0	0
5	B	1	Total Cl 1 1	0	0

- Molecule 6 is DINUCLEAR COPPER ION (CCD ID: CUA) (formula: Cu_2) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total Cu 2 2	0	0
6	B	1	Total Cu 2 2	0	0

- Molecule 7 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	B	1	Total Ca 1 1	0	0

- Molecule 8 is water.

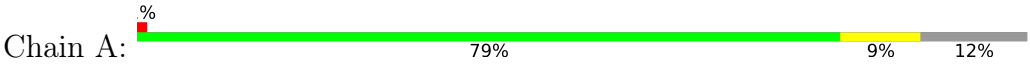
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	602	Total O 602 602	0	0
8	B	536	Total O 536 536	0	0

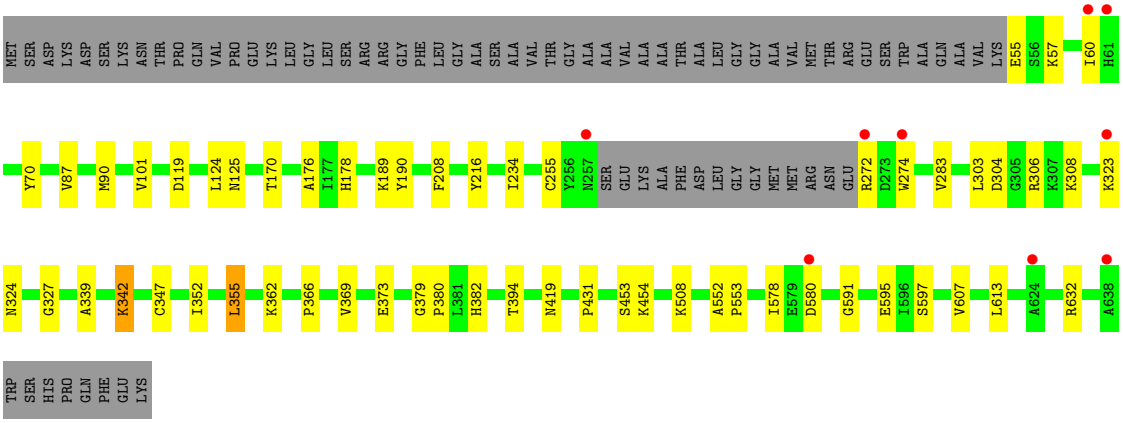
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Nitrous-oxide reductase

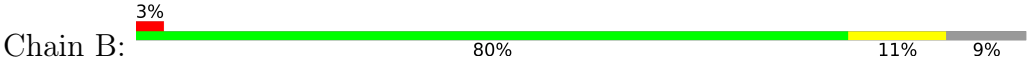
Chain A:

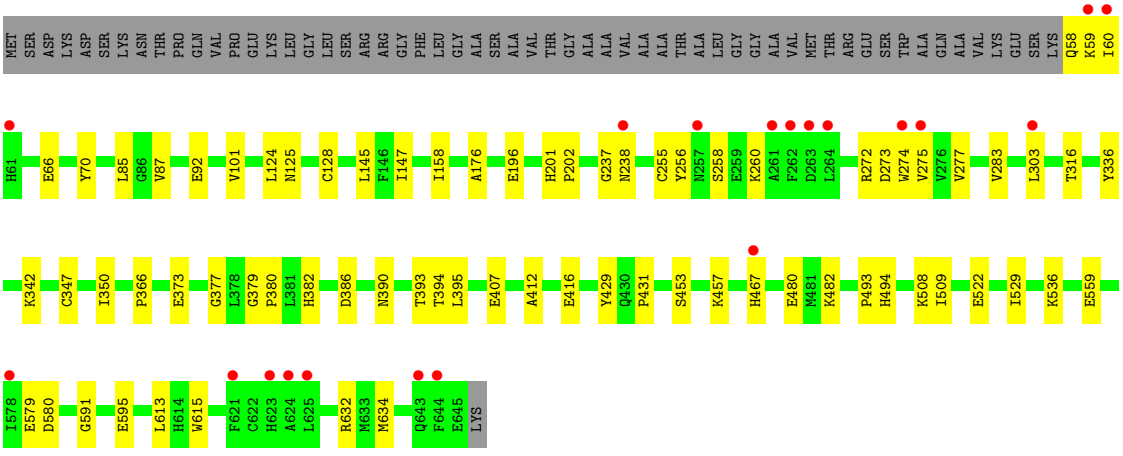




• Molecule 1: Nitrous-oxide reductase

Chain B:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	69.30Å 76.15Å 108.20Å 90.00° 93.19° 90.00°	Depositor
Resolution (Å)	76.15 – 1.60 76.15 – 1.60	Depositor EDS
% Data completeness (in resolution range)	74.8 (76.15-1.60) 67.8 (76.15-1.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.09 (at 1.59Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.160 , 0.200 0.162 , 0.201	Depositor DCC
R_{free} test set	5001 reflections (3.35%)	wwPDB-VP
Wilson B-factor (Å^2)	21.6	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 44.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	10347	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.17% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical 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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, FMT, CL, B3P, NA, CUA

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	0.34	0/4640	0.55	0/6285
1	B	0.31	0/4756	0.52	0/6443
All	All	0.32	0/9396	0.54	0/12728

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	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	336	TYR	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	4510	0	4401	44	0
1	B	4634	0	4497	57	0
2	A	12	0	4	0	0
2	B	6	0	2	2	0
3	A	19	0	26	0	0
3	B	19	0	26	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	2	0	0	0	0
6	B	2	0	0	0	0
7	B	1	0	0	0	0
8	A	602	0	0	18	0
8	B	536	0	0	11	1
All	All	10347	0	8956	98	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 5.

All (98) 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:304:ASP:OD2	8:A:801:HOH:O	1.94	0.84
1:A:352:ILE:HA	1:A:355:LEU:HD22	1.63	0.81
1:A:306:ARG:NE	8:A:801:HOH:O	2.15	0.78
3:B:902:B3P:HN2	2:B:903:FMT:H	1.51	0.74
1:B:522:GLU:OE1	8:B:1002:HOH:O	2.07	0.72
1:B:272:ARG:O	8:B:1003:HOH:O	2.08	0.71
1:B:508:LYS:NZ	1:B:580:ASP:OD1	2.21	0.69
1:B:595:GLU:OE2	8:B:1004:HOH:O	2.10	0.68
1:A:306:ARG:NH2	8:A:801:HOH:O	2.24	0.68
1:B:457:LYS:NZ	8:B:1007:HOH:O	2.28	0.66
1:B:416:GLU:OE2	8:B:1005:HOH:O	2.14	0.66
1:A:283:VAL:HG13	1:A:303:LEU:HD13	1.78	0.64
1:A:382:HIS:NE2	8:A:803:HOH:O	2.30	0.64
1:B:255[B]:CYS:HG	1:B:274:TRP:CD1	2.15	0.63
1:B:559:GLU:OE2	1:B:632:ARG:NH1	2.33	0.62
1:B:255[B]:CYS:SG	1:B:274:TRP:NE1	2.74	0.60
1:A:595:GLU:OE2	8:A:802:HOH:O	2.16	0.58
1:B:508:LYS:C	1:B:509:ILE:HD12	2.28	0.58
1:A:272:ARG:N	1:A:323:LYS:HZ2	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:362:LYS:NZ	8:A:815:HOH:O	2.36	0.57
1:A:308:LYS:NZ	8:A:801:HOH:O	2.33	0.56
1:A:101:VAL:HG21	1:A:124:LEU:HD22	1.87	0.56
1:B:632:ARG:NH1	8:B:1006:HOH:O	2.16	0.54
1:A:578:ILE:HD12	1:A:578:ILE:H	1.73	0.54
1:A:306:ARG:CZ	8:A:801:HOH:O	2.48	0.53
1:B:59:LYS:NZ	1:B:66:GLU:OE1	2.29	0.53
1:A:632:ARG:HD3	8:A:828:HOH:O	2.08	0.53
1:B:275:VAL:HG11	1:B:350:ILE:HD11	1.91	0.52
1:B:529:ILE:HD13	1:B:536:LYS:HE3	1.91	0.52
1:B:70:TYR:HB2	1:B:87:VAL:HB	1.91	0.52
1:A:431:PRO:HA	1:A:453:SER:HA	1.93	0.51
1:B:613:LEU:HD21	1:B:632:ARG:HG2	1.91	0.51
1:B:366:PRO:HG3	8:B:1396:HOH:O	2.11	0.50
1:A:454:LYS:HD2	8:A:808:HOH:O	2.12	0.50
1:A:55:GLU:N	8:A:821:HOH:O	2.44	0.49
1:A:255:CYS:HB2	1:A:274:TRP:CD1	2.47	0.49
1:B:407:GLU:HG2	8:B:1236:HOH:O	2.11	0.49
1:A:55:GLU:HB3	1:A:57:LYS:HE3	1.95	0.48
1:B:85:LEU:HD22	1:B:92:GLU:HA	1.95	0.48
1:A:170:THR:HG21	2:B:901:FMT:H	1.96	0.48
1:B:101:VAL:HG21	1:B:124:LEU:HD22	1.97	0.47
1:A:189:LYS:HE3	1:A:190:TYR:CZ	2.49	0.47
1:A:347:CYS:O	1:A:373:GLU:HA	2.13	0.47
1:A:419:ASN:OD1	8:A:804:HOH:O	2.20	0.47
1:B:128:CYS:HB3	1:B:147:ILE:HD11	1.97	0.47
1:A:70:TYR:HB2	1:A:87:VAL:HB	1.97	0.47
1:B:632:ARG:HD3	8:B:1006:HOH:O	2.14	0.46
1:B:615:TRP:CH2	1:B:632:ARG:HD2	2.51	0.46
1:A:125:ASN:OD1	1:B:591:GLY:HA2	2.16	0.46
1:B:382:HIS:O	1:B:393:THR:HA	2.16	0.46
1:A:90:MET:HE2	8:A:902:HOH:O	2.16	0.46
1:B:272:ARG:NH1	8:B:1030:HOH:O	2.49	0.46
1:B:237:GLY:H	1:B:260:LYS:NZ	2.14	0.45
1:B:494:HIS:CE1	8:B:1001:HOH:O	2.68	0.45
1:B:255[B]:CYS:HG	1:B:274:TRP:NE1	2.13	0.45
1:B:275:VAL:HG11	1:B:350:ILE:CD1	2.46	0.45
1:A:178:HIS:HB3	8:A:836:HOH:O	2.16	0.45
1:B:145:LEU:HD22	1:B:158:ILE:HD12	1.98	0.45
1:B:196:GLU:O	1:B:238:ASN:HB2	2.17	0.45
1:A:304:ASP:CG	8:A:801:HOH:O	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:377:GLY:HA3	1:B:395:LEU:HD13	2.00	0.44
1:B:508:LYS:HG3	1:B:579:GLU:OE1	2.18	0.44
1:B:431:PRO:HA	1:B:453:SER:HA	1.99	0.44
1:A:580:ASP:OD1	1:B:467:HIS:HB2	2.17	0.44
1:B:275:VAL:CG1	1:B:350:ILE:HD11	2.47	0.44
1:A:379:GLY:N	1:A:380:PRO:HD3	2.33	0.43
1:B:429:TYR:CD2	1:B:457:LYS:HD2	2.53	0.43
1:B:201:HIS:HA	1:B:202:PRO:HA	1.86	0.43
1:B:480:GLU:OE1	1:B:482:LYS:NZ	2.49	0.43
1:A:613:LEU:HD11	1:A:632:ARG:HG2	2.00	0.43
1:A:380:PRO:HA	1:A:394:THR:O	2.18	0.43
1:B:379:GLY:N	1:B:380:PRO:HD3	2.34	0.43
1:A:324:ASN:HB3	1:A:342:LYS:HZ2	1.83	0.42
1:B:347:CYS:O	1:B:373:GLU:HA	2.18	0.42
1:B:509:ILE:CD1	1:B:579:GLU:HG2	2.50	0.42
1:A:552:ALA:HA	1:A:553:PRO:HA	1.81	0.42
1:A:591:GLY:HA2	1:B:125:ASN:OD1	2.18	0.42
1:B:283:VAL:HG13	1:B:303:LEU:HD13	2.00	0.42
1:B:493:PRO:O	1:B:494:HIS:HB3	2.20	0.42
1:B:380:PRO:HA	1:B:394:THR:O	2.19	0.42
1:A:119:ASP:HB2	8:A:1247:HOH:O	2.19	0.42
1:A:216:TYR:CD1	1:A:234:ILE:HG23	2.54	0.42
1:B:509:ILE:HD11	1:B:579:GLU:HG2	2.01	0.42
1:A:208:PHE:CD1	1:B:634:MET:HG3	2.55	0.41
1:A:60:ILE:HG22	8:A:966:HOH:O	2.21	0.41
1:B:58:GLN:O	1:B:60:ILE:HG13	2.20	0.41
1:B:386:ASP:OD2	1:B:390:ASN:HB2	2.20	0.41
1:B:634:MET:HE2	1:B:634:MET:HB2	1.92	0.41
1:B:238:ASN:ND2	1:B:256:TYR:CZ	2.85	0.41
1:B:412:ALA:HA	1:B:416:GLU:O	2.20	0.41
1:A:366:PRO:O	1:A:369:VAL:HG22	2.19	0.41
1:A:324:ASN:HB3	1:A:342:LYS:NZ	2.36	0.40
1:B:258:SER:OG	1:B:273:ASP:OD2	2.29	0.40
1:B:277:VAL:O	1:B:316:THR:HA	2.21	0.40
1:A:508:LYS:HD2	8:A:856:HOH:O	2.21	0.40
1:A:595:GLU:OE1	1:A:597:SER:OG	2.36	0.40
1:B:255[B]:CYS:SG	1:B:274:TRP:CD1	3.12	0.40
1:A:327:GLY:O	1:A:339:ALA:HA	2.22	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)
8:B:1250:HOH:O	8:B:1450:HOH:O[1_565]	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	573/646 (89%)	546 (95%)	25 (4%)	2 (0%)	36 20
1	B	588/646 (91%)	564 (96%)	22 (4%)	2 (0%)	36 20
All	All	1161/1292 (90%)	1110 (96%)	47 (4%)	4 (0%)	36 20

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	342	LYS
1	B	342	LYS
1	A	176	ALA
1	B	176	ALA

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	494/544 (91%)	492 (100%)	2 (0%)	84 75
1	B	504/544 (93%)	504 (100%)	0	100 100
All	All	998/1088 (92%)	996 (100%)	2 (0%)	87 81

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	355	LEU
1	A	607	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	HIS
1	A	111	ASN
1	A	178	HIS
1	A	577	GLN
1	A	599	GLN
1	B	232	GLN
1	B	390	ASN
1	B	467	HIS
1	B	530	ASN
1	B	599	GLN
1	B	643	GLN

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

Of 15 ligands modelled in this entry, 5 are monoatomic - leaving 10 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	FMT	A	704	1	2,2,2	0.58	0	1,1,1	0.13	0
2	FMT	A	702	-	2,2,2	0.70	0	1,1,1	0.24	0
2	FMT	A	703	-	2,2,2	0.69	0	1,1,1	0.06	0
3	B3P	B	902	-	18,18,18	0.77	0	21,23,23	1.06	1 (4%)
2	FMT	B	901	-	2,2,2	0.63	0	1,1,1	0.07	0
6	CUA	B	907	1	0,1,1	-	-	-		
3	B3P	A	705	-	18,18,18	0.69	0	21,23,23	0.91	1 (4%)
2	FMT	B	903	-	2,2,2	0.68	0	1,1,1	0.12	0
2	FMT	A	701	-	2,2,2	0.63	0	1,1,1	0.08	0
6	CUA	A	708	1	0,1,1	-	-	-		

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	B3P	B	902	-	-	5/28/28/28	-
3	B3P	A	705	-	-	0/28/28/28	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	902	B3P	O2-C10-C8	-2.06	107.46	111.63
3	A	705	B3P	O5-C6-C4	-2.02	107.54	111.63

There are no chirality outliers.

All (5) torsion outliers are listed below:

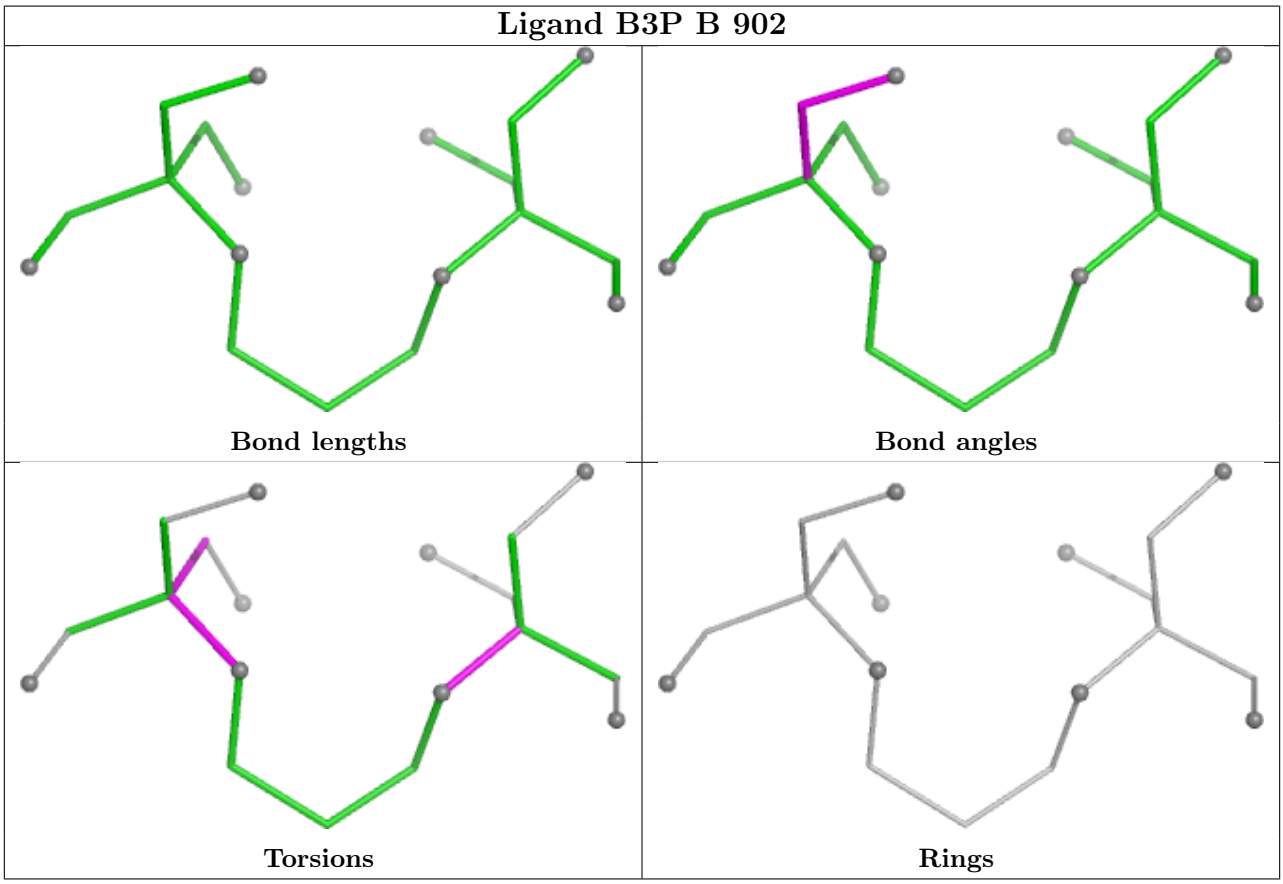
Mol	Chain	Res	Type	Atoms
3	B	902	B3P	C5-C4-N1-C3
3	B	902	B3P	C11-C8-N2-C2
3	B	902	B3P	C11-C8-C9-O1
3	B	902	B3P	C10-C8-N2-C2
3	B	902	B3P	C9-C8-N2-C2

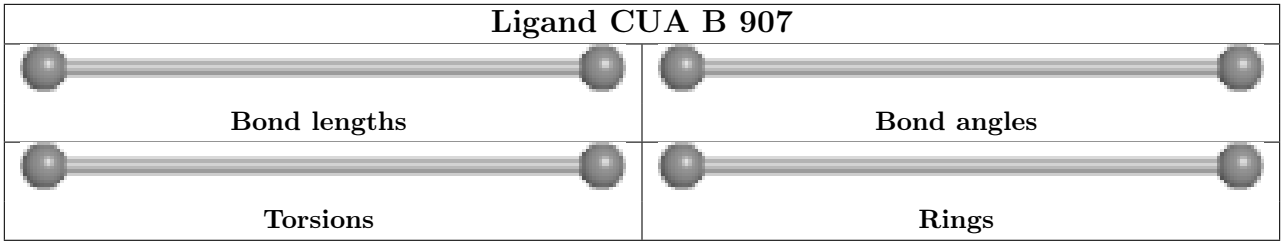
There are no ring outliers.

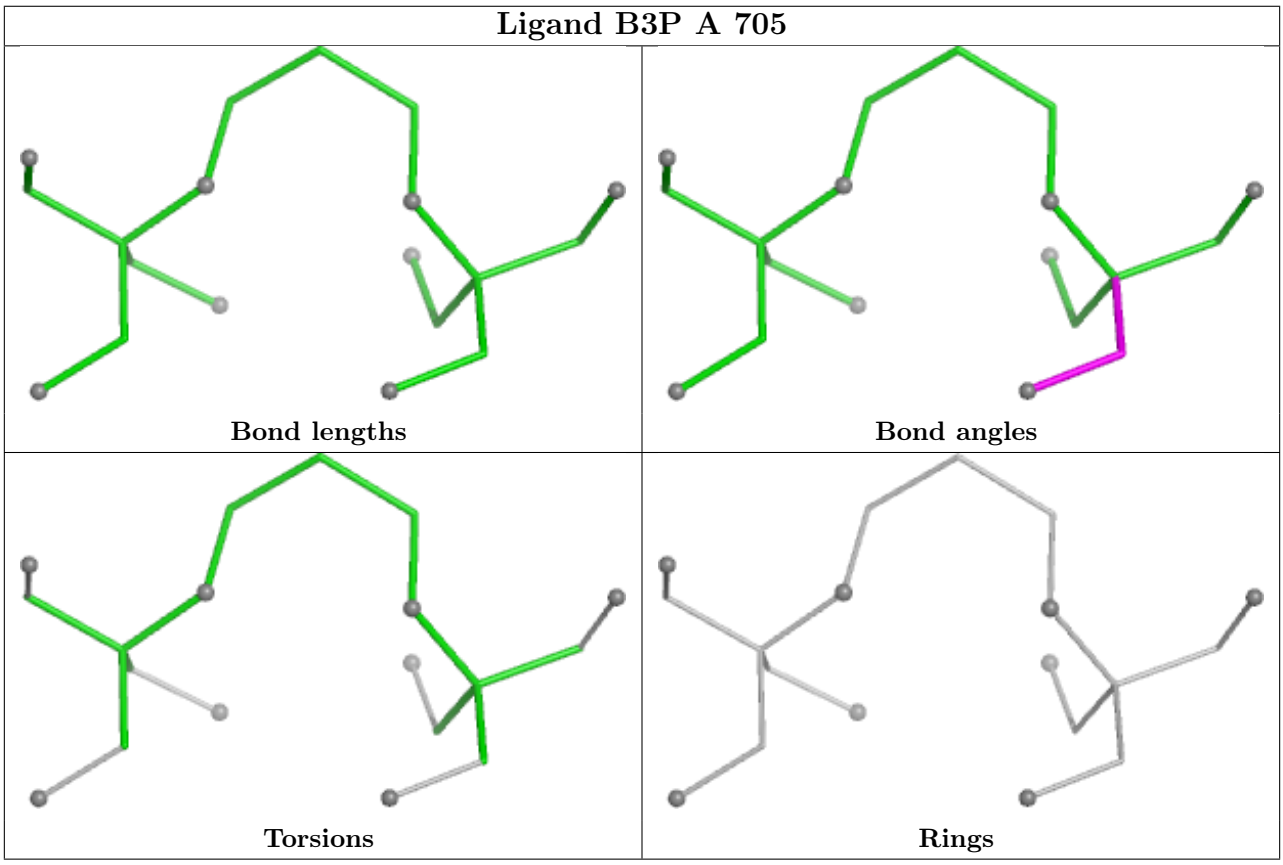
3 monomers are involved in 2 short contacts:

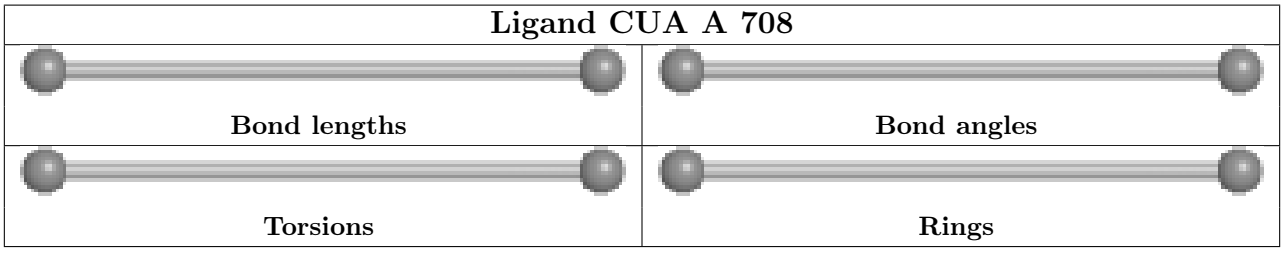
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	902	B3P	1	0
2	B	901	FMT	1	0
2	B	903	FMT	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	570/646 (88%)	-0.36	9 (1%) 70 74	9, 23, 46, 72	7 (1%)
1	B	588/646 (91%)	-0.10	20 (3%) 48 50	12, 28, 56, 80	2 (0%)
All	All	1158/1292 (89%)	-0.23	29 (2%) 58 61	9, 25, 51, 80	9 (0%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	638	ALA	5.5
1	B	60	ILE	3.6
1	B	621	PHE	3.4
1	B	625	LEU	2.9
1	B	262	PHE	2.9
1	A	257	ASN	2.9
1	B	264	LEU	2.9
1	A	60	ILE	2.9
1	B	61	HIS	2.8
1	B	644	PHE	2.8
1	B	303	LEU	2.8
1	B	261	ALA	2.7
1	B	624	ALA	2.5
1	A	323	LYS	2.5
1	A	272	ARG	2.3
1	B	578	ILE	2.3
1	B	238	ASN	2.3
1	B	263	ASP	2.3
1	B	257	ASN	2.2
1	A	61	HIS	2.2
1	A	274	TRP	2.2
1	B	623	HIS	2.1
1	B	274	TRP	2.1
1	B	275	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	643	GLN	2.1
1	A	624	ALA	2.1
1	B	467	HIS	2.0
1	B	59	LYS	2.0
1	A	580	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, 95^{th} 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.

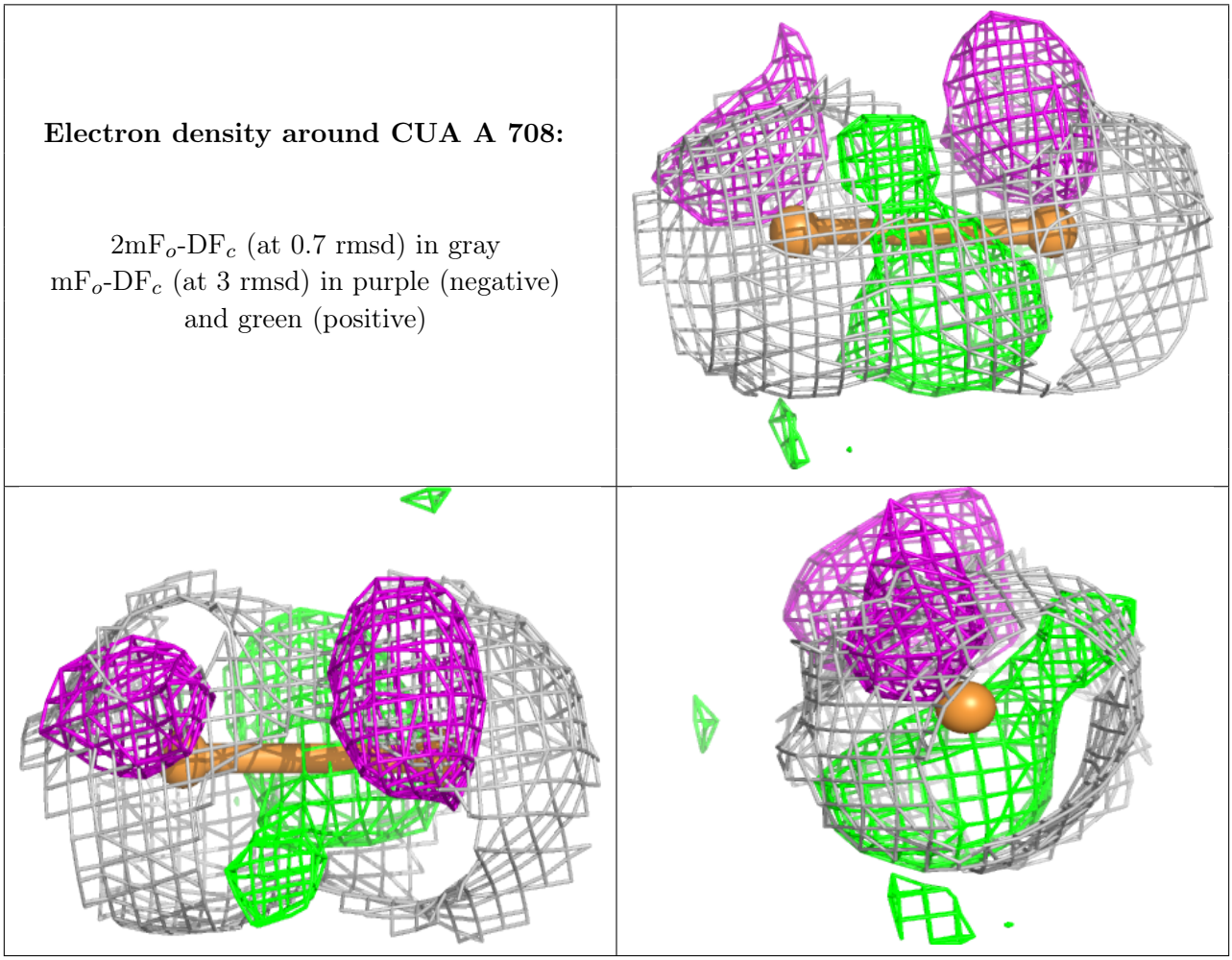
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	FMT	A	701	3/3	0.78	0.15	31,31,46,52	0
2	FMT	A	704	3/3	0.78	0.15	31,31,45,57	0
2	FMT	B	901	3/3	0.85	0.25	35,35,37,39	0
2	FMT	A	703	3/3	0.91	0.20	37,37,41,49	0
6	CUA	A	708	2/2	0.92	0.08	40,40,40,45	2
6	CUA	B	907	2/2	0.94	0.07	37,37,37,47	1
3	B3P	A	705	19/19	0.95	0.06	16,23,29,30	0
2	FMT	B	903	3/3	0.96	0.12	19,19,33,40	0
2	FMT	A	702	3/3	0.96	0.13	21,21,38,44	0
3	B3P	B	902	19/19	0.97	0.05	18,23,27,30	0
5	CL	B	906	1/1	0.98	0.04	27,27,27,27	1
7	CA	B	905	1/1	0.98	0.04	24,24,24,24	1
4	NA	B	904	1/1	0.99	0.03	24,24,24,24	1
5	CL	A	707	1/1	0.99	0.04	21,21,21,21	1
4	NA	A	706	1/1	0.99	0.03	19,19,19,19	1

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.

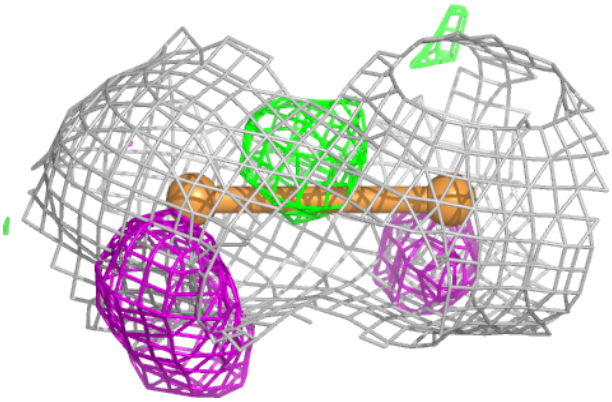
Electron density around CUA A 708:

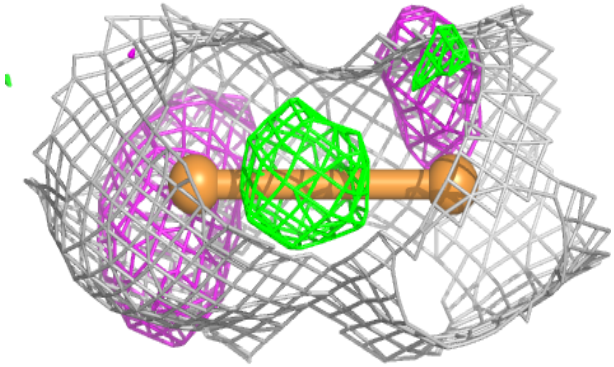
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

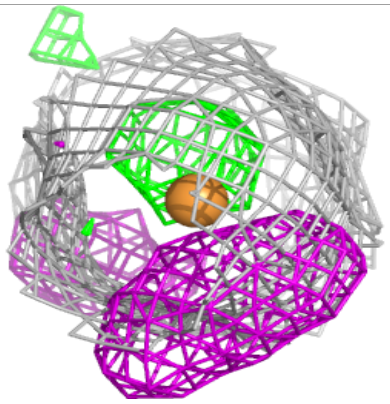


Electron density around CUA B 907:

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

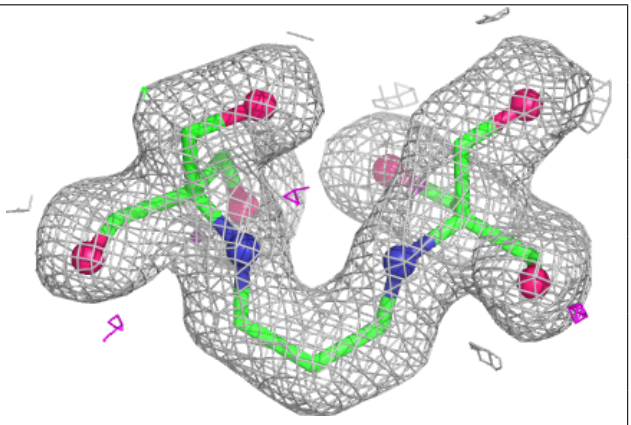


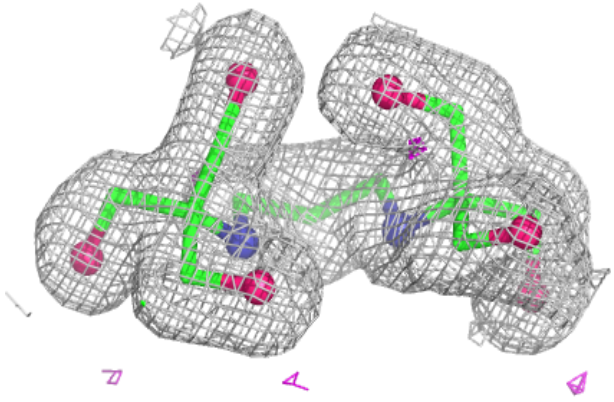


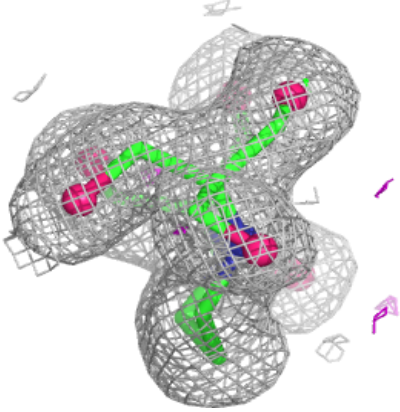


Electron density around B3P A 705:

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

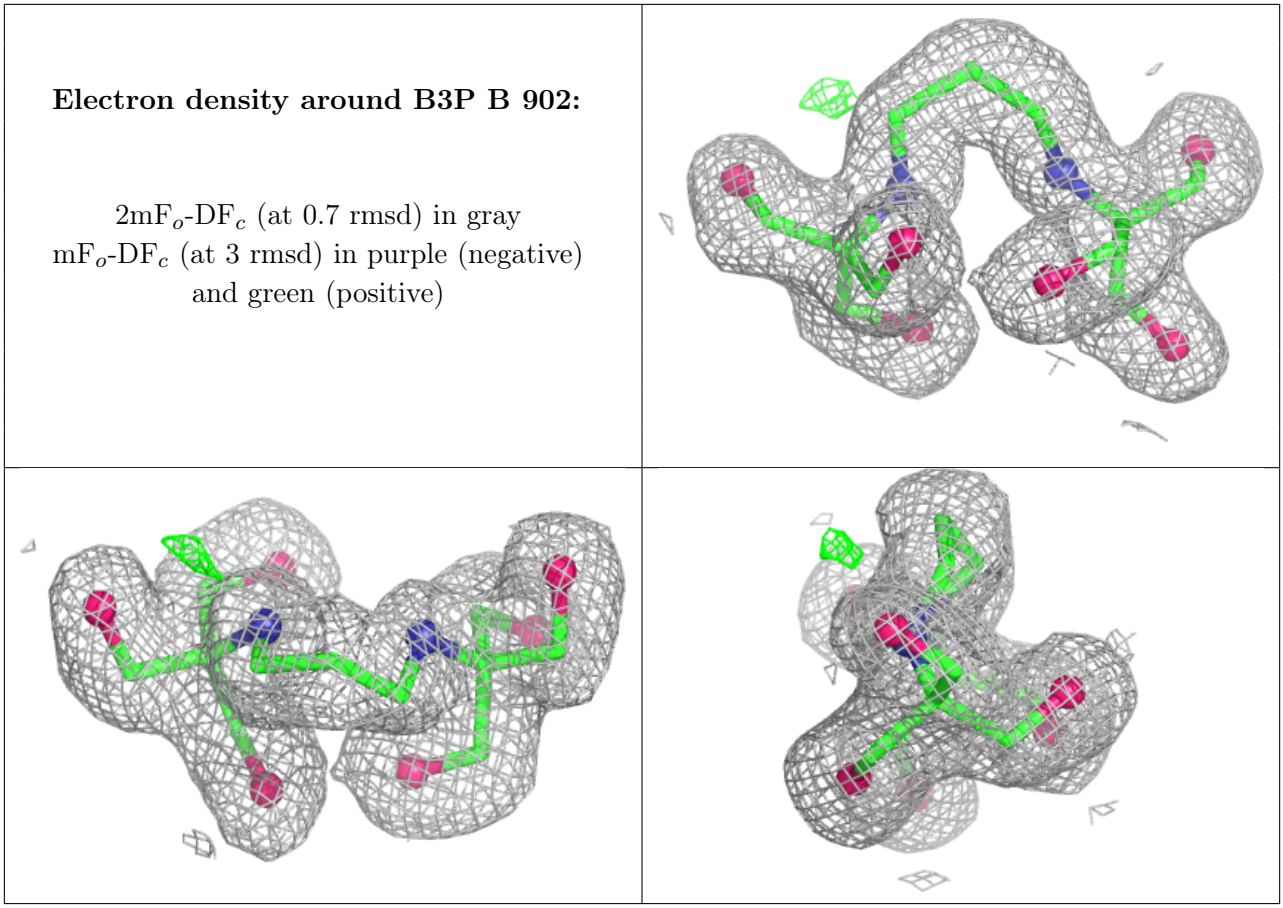






Electron density around B3P B 902:

$2mF_o$-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.