



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 05:58 PM UTC

PDB ID : 8CGE / pdb_00008cge
Title : The crystal structure of a cobalt-bound scFv reveals a Tetrameric polyHistidine motif (TetrHis)
Authors : Healey, R.D.; Hoh, F.; Granier, S.; Leyrat, C.
Deposited on : 2023-02-03
Resolution : 2.50 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
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

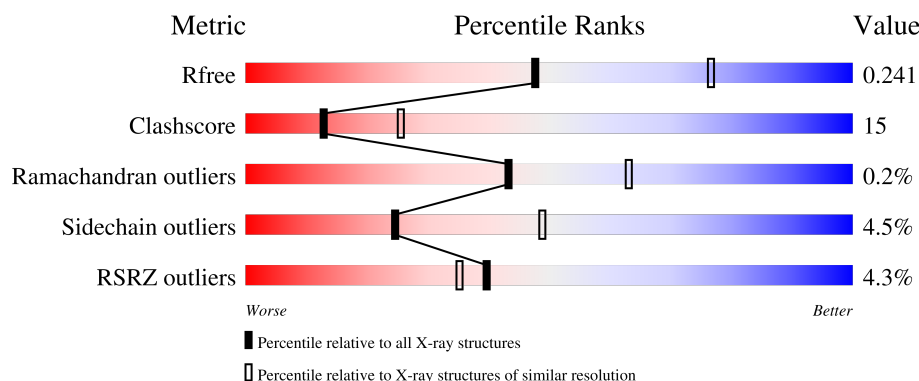
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	304	5% 60% 19% • 19%
1	B	304	2% 57% 22% • 19%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4154 atoms, of which 19 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.

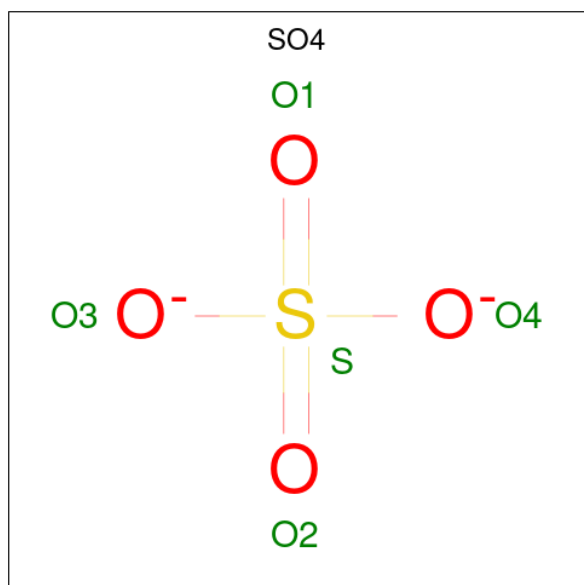
- Molecule 1 is a protein called Single Chain Fv 2A2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total	C	N	O	S	0	0	0
			1913	1212	321	373	7			
1	B	246	Total	C	N	O	S	0	0	0
			1933	1227	323	376	7			

- Molecule 2 is COBALT (II) ION (CCD ID: CO) (formula: Co) (labeled as "Ligand of Interest" by depositor).

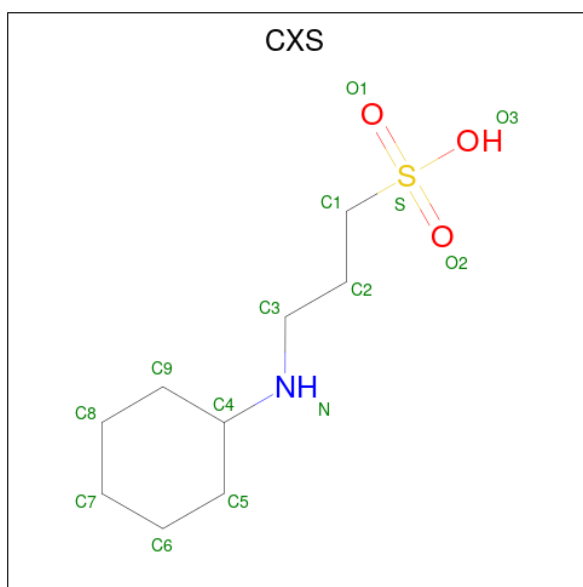
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Co	0	0
			2	2		
2	B	2	Total	Co	0	0
			2	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0

- Molecule 4 is 3-CYCLOHEXYL-1-PROPYLSULFONIC ACID (CCD ID: CXS) (formula: $C_9H_{19}NO_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	B	1	Total	C	H	N	O	S	25	0
			33	9	19	1	3	1		

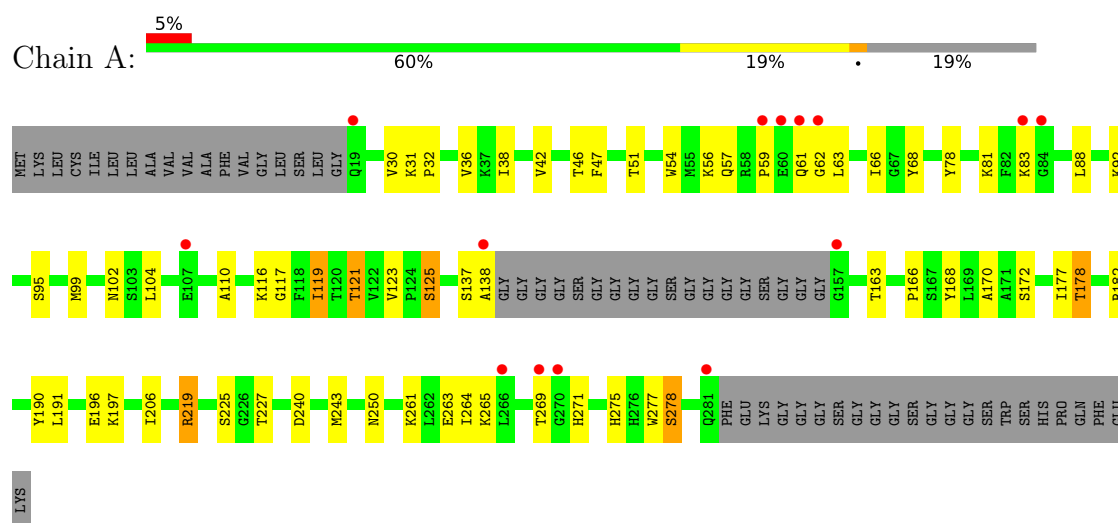
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	101	Total	O	0	0
			101	101		
5	B	95	Total	O	0	0
			95	95		

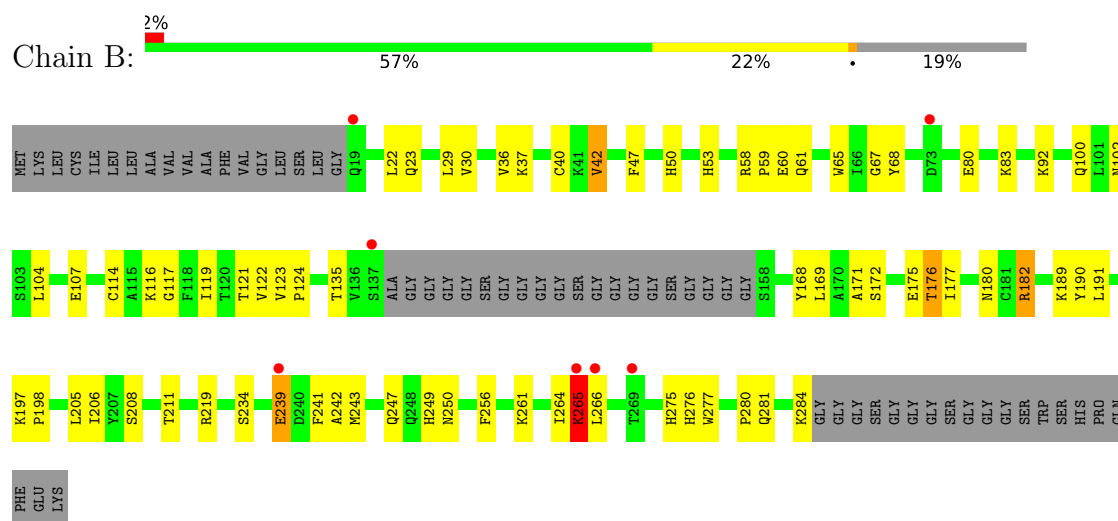
3 Residue-property plots [i](#)

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: Single Chain Fv 2A2



• Molecule 1: Single Chain Fv 2A2



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	97.50Å 81.82Å 111.05Å 90.00° 111.14° 90.00°	Depositor
Resolution (Å)	45.17 – 2.50 45.17 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.0 (45.17-2.50) 99.2 (45.17-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.51Å)	Xtriage
Refinement program	BUSTER, PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.190 , 0.248 (Not available) , 0.241	Depositor DCC
R_{free} test set	1389 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	44.8	Xtriage
Anisotropy	0.314	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4154	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.61% 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: CO, CXS, SO4

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	0.37	0/1967	0.67	5/2670 (0.2%)
1	B	0.35	0/1988	0.66	6/2697 (0.2%)
All	All	0.36	0/3955	0.66	11/5367 (0.2%)

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.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	83	LYS	CG-CD-CE	10.95	136.49	111.30
1	A	83	LYS	CB-CG-CD	-7.46	94.15	111.30
1	B	239	GLU	CA-CB-CG	6.83	127.76	114.10
1	B	265	LYS	CA-C-N	-6.32	100.23	122.15
1	B	265	LYS	C-N-CA	-6.32	100.23	122.15
1	B	265	LYS	CA-CB-CG	5.71	125.52	114.10
1	B	265	LYS	CB-CA-C	-5.69	100.27	109.72
1	A	81	LYS	CA-CB-CG	5.46	125.02	114.10
1	A	83	LYS	N-CA-CB	-5.31	102.16	109.97
1	A	81	LYS	CD-CE-NZ	5.19	128.51	111.90
1	B	239	GLU	CB-CA-C	-5.04	99.78	110.31

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	265	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	1913	0	1828	51	0
1	B	1933	0	1848	73	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	40	0	0	1	0
3	B	35	0	0	0	0
4	B	14	19	18	0	0
5	A	101	0	0	10	0
5	B	95	0	0	3	0
All	All	4135	19	3694	115	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 15.

All (115) 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:166:PRO:HG3	1:B:169:LEU:HD11	1.32	1.09
1:B:265:LYS:O	1:B:266:LEU:HD22	1.61	1.00
1:B:172:SER:HB2	1:B:175:GLU:CD	1.99	0.87
1:A:178:THR:HG22	1:B:180:ASN:HD21	1.37	0.87
1:A:197:LYS:NZ	5:A:501:HOH:O	2.10	0.85
1:A:166:PRO:CG	1:B:169:LEU:HD11	2.05	0.85
1:B:30:VAL:HG11	1:B:36:VAL:CG1	2.10	0.82
1:B:30:VAL:HG21	1:B:104:LEU:CD1	2.13	0.79
1:B:172:SER:OG	1:B:265:LYS:HE2	1.83	0.77
1:B:168:TYR:CD1	1:B:261:LYS:HE2	2.21	0.76
1:B:30:VAL:HG21	1:B:104:LEU:HD12	1.67	0.76
1:B:59:PRO:HD3	1:B:281:GLN:NE2	2.03	0.72
1:A:92:LYS:NZ	5:A:504:HOH:O	2.22	0.72
1:A:51:THR:HG23	1:A:119:ILE:CD1	2.20	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:THR:HG23	1:A:271:HIS:H	1.56	0.70
1:B:172:SER:HA	1:B:265:LYS:HG3	1.72	0.69
1:B:197:LYS:HG2	1:B:242:ALA:HB2	1.75	0.68
1:B:59:PRO:HD3	1:B:281:GLN:HE21	1.58	0.67
1:B:172:SER:OG	1:B:265:LYS:CE	2.43	0.67
1:A:178:THR:CG2	1:B:180:ASN:HD21	2.08	0.66
1:B:30:VAL:HG11	1:B:36:VAL:HG13	1.78	0.66
1:B:175:GLU:OE2	5:B:501:HOH:O	2.14	0.66
1:A:62:GLY:HA2	5:A:503:HOH:O	1.95	0.65
1:B:172:SER:CA	1:B:265:LYS:HG3	2.27	0.65
1:B:168:TYR:HD1	1:B:261:LYS:HE2	1.61	0.65
1:A:51:THR:HG23	1:A:119:ILE:HD11	1.79	0.64
1:B:36:VAL:HG22	1:B:104:LEU:HD11	1.80	0.64
1:B:172:SER:HB2	1:B:175:GLU:CG	2.28	0.63
1:B:247:GLN:HG3	1:B:256:PHE:CE2	2.33	0.63
1:B:172:SER:OG	1:B:265:LYS:HG3	1.98	0.63
1:A:32:PRO:HD3	1:A:137:SER:O	1.98	0.63
1:B:182:ARG:HB2	1:B:182:ARG:HH11	1.64	0.62
1:A:275:HIS:HE1	5:A:566:HOH:O	1.82	0.62
1:B:42:VAL:HG21	1:B:47:PHE:HD1	1.65	0.62
1:A:36:VAL:HG22	1:A:104:LEU:HD11	1.80	0.61
1:A:121:THR:HB	1:A:123:VAL:H	1.65	0.60
1:A:168:TYR:HE2	1:A:263:GLU:OE1	1.86	0.59
1:B:42:VAL:HG21	1:B:47:PHE:CD1	2.39	0.58
1:A:56:LYS:NZ	5:A:509:HOH:O	2.36	0.57
1:B:29:LEU:HD13	1:B:135:THR:HB	1.86	0.57
1:B:65:TRP:CZ2	1:B:67:GLY:HA2	2.40	0.57
1:A:63:LEU:N	5:A:503:HOH:O	2.11	0.56
1:B:37:LYS:NZ	5:B:502:HOH:O	2.39	0.56
1:B:22:LEU:HD23	1:B:40:CYS:SG	2.46	0.55
1:A:170:ALA:HB1	1:A:265:LYS:HA	1.88	0.55
1:A:243:MET:HE1	1:A:261:LYS:HD2	1.88	0.55
1:B:60:GLU:HB3	1:B:61:GLN:OE1	2.07	0.54
1:B:265:LYS:C	1:B:266:LEU:HD22	2.30	0.54
1:B:172:SER:HB2	1:B:175:GLU:HG3	1.90	0.54
1:B:172:SER:O	1:B:175:GLU:HG3	2.08	0.54
1:B:58:ARG:NH1	1:B:107:GLU:OE2	2.35	0.53
1:B:30:VAL:HG11	1:B:36:VAL:HG11	1.87	0.53
1:A:30:VAL:HG21	1:A:104:LEU:HD13	1.91	0.53
1:A:102:ASN:ND2	5:A:502:HOH:O	2.10	0.53
1:A:172:SER:OG	1:A:265:LYS:HD3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:219:ARG:NH2	1:B:239:GLU:OE2	2.43	0.51
1:A:42:VAL:HG21	1:A:47:PHE:HD1	1.75	0.51
1:A:166:PRO:CB	1:B:169:LEU:HD11	2.40	0.51
1:B:169:LEU:HD23	1:B:177:ILE:HD13	1.93	0.51
1:B:80:GLU:OE1	1:B:83:LYS:NZ	2.39	0.50
1:A:54:TRP:CD1	1:A:88:LEU:HD22	2.46	0.50
5:A:551:HOH:O	1:B:176:THR:HG22	2.09	0.50
1:B:172:SER:OG	1:B:265:LYS:CD	2.60	0.50
1:A:119:ILE:HG23	1:A:119:ILE:O	2.11	0.50
1:A:59:PRO:O	1:A:61:GLN:NE2	2.28	0.49
1:B:172:SER:CB	1:B:265:LYS:HG3	2.42	0.49
1:A:182:ARG:HA	1:A:227:THR:O	2.12	0.48
1:A:116:LYS:O	1:A:125:SER:HB3	2.13	0.48
1:B:53:HIS:O	1:B:114:CYS:HA	2.14	0.47
1:A:54:TRP:CE3	1:A:99:MET:HE2	2.50	0.47
1:A:219:ARG:NH1	1:A:240:ASP:OD1	2.48	0.47
1:B:172:SER:OG	1:B:265:LYS:CG	2.61	0.47
1:A:278:SER:HB2	1:B:266:LEU:HD23	1.96	0.47
1:B:176:THR:HB	1:B:234:SER:HA	1.97	0.46
1:A:178:THR:HG22	1:B:180:ASN:ND2	2.18	0.46
1:A:57:GLN:HB2	1:A:63:LEU:HD23	1.98	0.46
1:B:121:THR:OG1	1:B:123:VAL:HG23	2.16	0.45
1:A:31:LYS:HG2	1:A:138:ALA:O	2.17	0.45
1:A:51:THR:OG1	1:A:117:GLY:HA3	2.17	0.45
1:A:57:GLN:O	1:A:110:ALA:HB1	2.16	0.45
1:A:219:ARG:NH1	1:A:240:ASP:OD2	2.49	0.45
1:B:190:TYR:HB2	1:B:250:ASN:HB2	1.98	0.45
1:B:50:HIS:CD2	1:B:116:LYS:HE2	2.52	0.45
1:A:190:TYR:HB2	1:A:250:ASN:HB2	1.99	0.45
1:B:50:HIS:HB3	1:B:117:GLY:O	2.16	0.45
1:A:277:TRP:CE2	1:B:275:HIS:HB2	2.53	0.44
1:B:68:TYR:CD1	1:B:68:TYR:C	2.95	0.44
1:B:198:PRO:HG3	1:B:277:TRP:HB2	1.98	0.44
1:A:177:ILE:C	1:A:177:ILE:HD12	2.43	0.44
1:B:23:GLN:NE2	1:B:23:GLN:HA	2.32	0.43
1:B:243:MET:HE1	1:B:261:LYS:HB2	2.00	0.43
1:B:197:LYS:HE3	1:B:241:PHE:O	2.19	0.43
1:A:42:VAL:HG21	1:A:47:PHE:CD1	2.54	0.43
1:A:56:LYS:HB2	1:A:66:ILE:HD11	2.00	0.42
1:B:171:ALA:O	1:B:264:ILE:HA	2.19	0.42
1:A:38:ILE:HB	5:A:554:HOH:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:169:LEU:HD12	1:B:169:LEU:HA	1.66	0.42
1:A:78:TYR:HD1	3:A:410:SO4:O3	2.02	0.42
1:B:205:LEU:C	1:B:206:ILE:HG12	2.44	0.42
1:A:261:LYS:HG2	1:A:263:GLU:HG3	2.02	0.42
5:A:598:HOH:O	1:B:176:THR:HG21	2.19	0.42
1:B:122:VAL:O	1:B:122:VAL:HG12	2.20	0.42
1:B:169:LEU:HD21	1:B:177:ILE:HB	2.00	0.42
1:A:68:TYR:CD1	1:A:68:TYR:C	2.98	0.41
1:B:124:PRO:HD3	1:B:249:HIS:CG	2.55	0.41
1:B:247:GLN:HG3	1:B:256:PHE:CD2	2.55	0.41
1:B:30:VAL:HG21	1:B:104:LEU:HD13	1.96	0.41
1:A:275:HIS:O	1:B:276:HIS:HA	2.21	0.41
1:A:170:ALA:HA	1:A:263:GLU:O	2.20	0.41
1:B:189:LYS:O	1:B:208:SER:C	2.64	0.41
1:B:36:VAL:O	1:B:100:GLN:HA	2.21	0.41
1:B:280:PRO:O	1:B:284:LYS:HD2	2.20	0.41
1:B:92:LYS:HE2	5:B:584:HOH:O	2.21	0.40
1:A:31:LYS:HB2	1:A:31:LYS:HE3	1.73	0.40
1:A:47:PHE:CD2	1:A:95:SER:HA	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	241/304 (79%)	234 (97%)	7 (3%)	0	100	100
1	B	242/304 (80%)	237 (98%)	4 (2%)	1 (0%)	30	49
All	All	483/608 (79%)	471 (98%)	11 (2%)	1 (0%)	43	63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	119	ILE

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	210/241 (87%)	197 (94%)	13 (6%)	16	34
1	B	213/241 (88%)	207 (97%)	6 (3%)	38	66
All	All	423/482 (88%)	404 (96%)	19 (4%)	24	49

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	THR
1	A	119	ILE
1	A	121	THR
1	A	125	SER
1	A	163	THR
1	A	178	THR
1	A	191	LEU
1	A	196	GLU
1	A	206	ILE
1	A	219	ARG
1	A	225	SER
1	A	264	ILE
1	A	278	SER
1	B	42	VAL
1	B	102	ASN
1	B	176	THR
1	B	182	ARG
1	B	191	LEU
1	B	211	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	57	GLN
1	A	250	ASN
1	B	23	GLN
1	B	180	ASN
1	B	276	HIS
1	B	281	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 20 ligands modelled in this entry, 4 are monoatomic - leaving 16 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	$\# Z > 2$
3	SO4	A	410	-	4,4,4	0.25	0	6,6,6	0.15	0
3	SO4	B	410	-	4,4,4	0.27	0	6,6,6	0.21	0
3	SO4	B	408	-	4,4,4	0.27	0	6,6,6	0.11	0
3	SO4	A	407	-	4,4,4	0.27	0	6,6,6	0.09	0
3	SO4	B	404	-	4,4,4	0.28	0	6,6,6	0.12	0
3	SO4	A	403	-	4,4,4	0.29	0	6,6,6	0.15	0
3	SO4	B	406	-	4,4,4	0.25	0	6,6,6	0.11	0
3	SO4	B	407	-	4,4,4	0.18	0	6,6,6	0.24	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	B	409	-	4,4,4	0.29	0	6,6,6	0.17	0
3	SO4	A	405	-	4,4,4	0.24	0	6,6,6	0.13	0
4	CXS	B	401	-	14,14,14	0.55	0	18,18,18	1.94	4 (22%)
3	SO4	A	408	-	4,4,4	0.29	0	6,6,6	0.20	0
3	SO4	A	406	-	4,4,4	0.25	0	6,6,6	0.17	0
3	SO4	A	404	-	4,4,4	0.26	0	6,6,6	0.18	0
3	SO4	A	409	-	4,4,4	0.24	0	6,6,6	0.07	0
3	SO4	B	405	-	4,4,4	0.26	0	6,6,6	0.17	0

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
4	CXS	B	401	-	-	0/8/16/16	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	401	CXS	C5-C4-N	4.66	133.84	110.56
4	B	401	CXS	C3-N-C4	4.62	123.07	114.18
4	B	401	CXS	C9-C4-N	-3.38	93.68	110.56
4	B	401	CXS	C2-C1-S	-2.99	108.66	113.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	410	SO4	1	0

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, 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	245/304 (80%)	0.23	14 (5%) 29 26	31, 44, 71, 121	0
1	B	246/304 (80%)	0.05	7 (2%) 55 50	34, 43, 64, 97	0
All	All	491/608 (80%)	0.14	21 (4%) 40 35	31, 44, 68, 121	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	265	LYS	5.8
1	A	61	GLN	5.6
1	A	59	PRO	5.2
1	B	239	GLU	3.9
1	A	107	GLU	3.4
1	A	138	ALA	3.2
1	A	19	GLN	3.1
1	B	19	GLN	3.0
1	B	73	ASP	2.9
1	A	60	GLU	2.8
1	A	281	GLN	2.6
1	A	269	THR	2.5
1	A	83	LYS	2.5
1	A	266	LEU	2.4
1	A	84	GLY	2.3
1	A	270	GLY	2.3
1	A	157	GLY	2.2
1	A	62	GLY	2.2
1	B	269	THR	2.1
1	B	137	SER	2.1
1	B	266	LEU	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

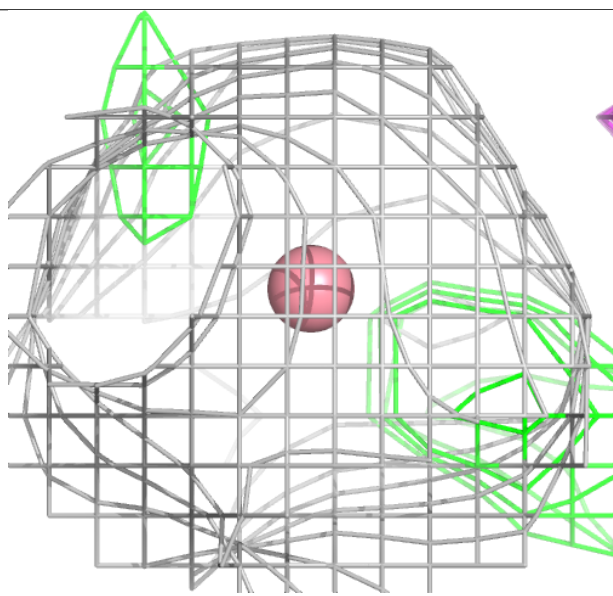
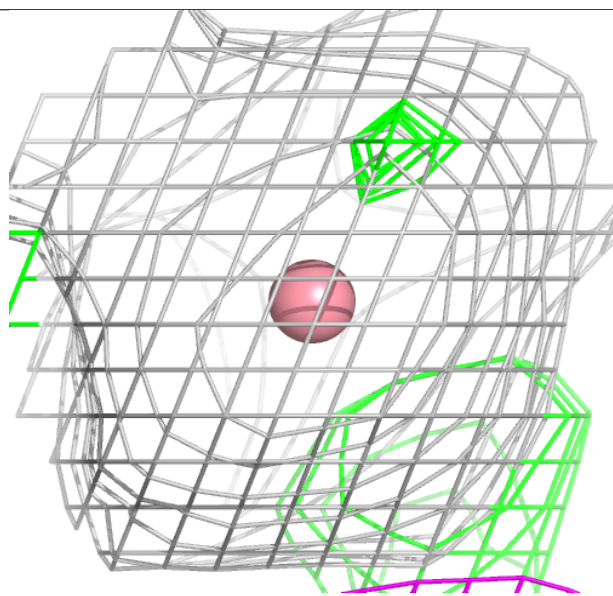
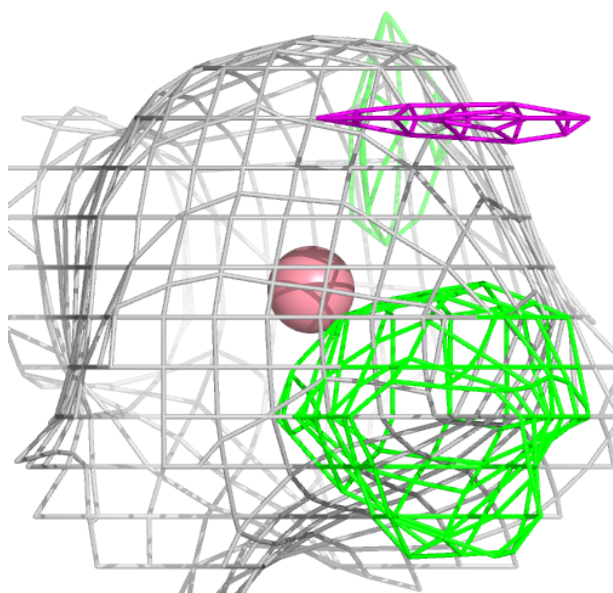
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	A	404	5/5	0.67	0.16	79,85,126,135	0
3	SO4	B	405	5/5	0.76	0.13	94,99,127,129	0
4	CXS	B	401	14/14	0.80	0.22	53,70,77,92	25
3	SO4	A	405	5/5	0.83	0.16	69,76,83,88	5
3	SO4	A	408	5/5	0.84	0.14	45,45,73,76	5
3	SO4	B	409	5/5	0.85	0.16	67,67,80,85	5
3	SO4	B	410	5/5	0.85	0.15	59,69,77,78	5
3	SO4	A	407	5/5	0.85	0.14	62,76,79,87	5
3	SO4	B	404	5/5	0.88	0.20	49,65,69,73	5
3	SO4	B	408	5/5	0.89	0.10	62,67,83,93	5
3	SO4	B	406	5/5	0.89	0.15	79,79,86,89	5
3	SO4	A	409	5/5	0.91	0.09	58,58,78,78	5
3	SO4	A	406	5/5	0.91	0.10	56,60,73,74	5
3	SO4	A	403	5/5	0.92	0.17	57,58,81,83	0
3	SO4	B	407	5/5	0.94	0.12	54,57,64,75	0
3	SO4	A	410	5/5	0.95	0.11	50,52,60,64	5
2	CO	A	402	1/1	0.97	0.06	67,67,67,67	0
2	CO	B	402	1/1	0.99	0.05	62,62,62,62	0
2	CO	B	403	1/1	0.99	0.02	51,51,51,51	0
2	CO	A	401	1/1	0.99	0.02	60,60,60,60	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.

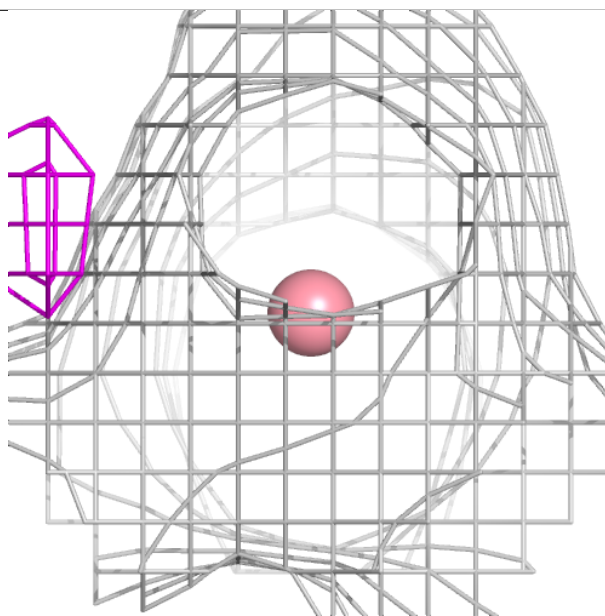
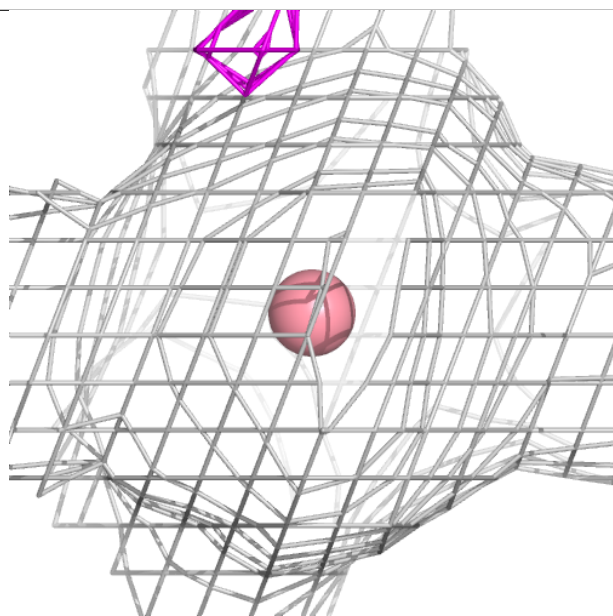
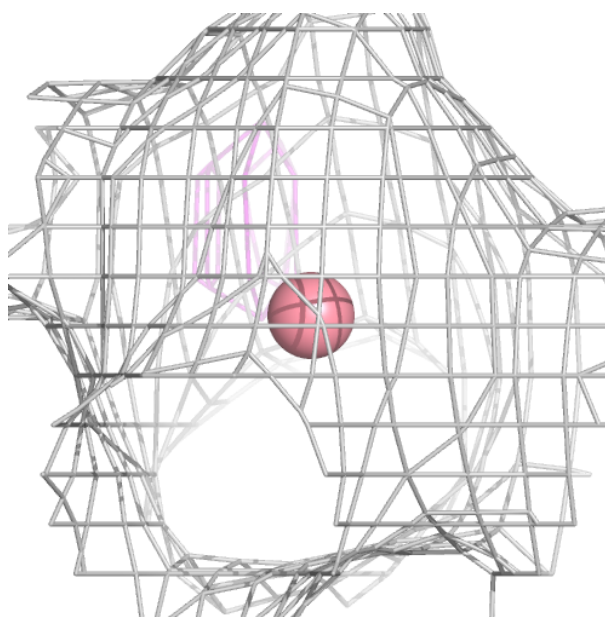
Electron density around CO A 402:

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



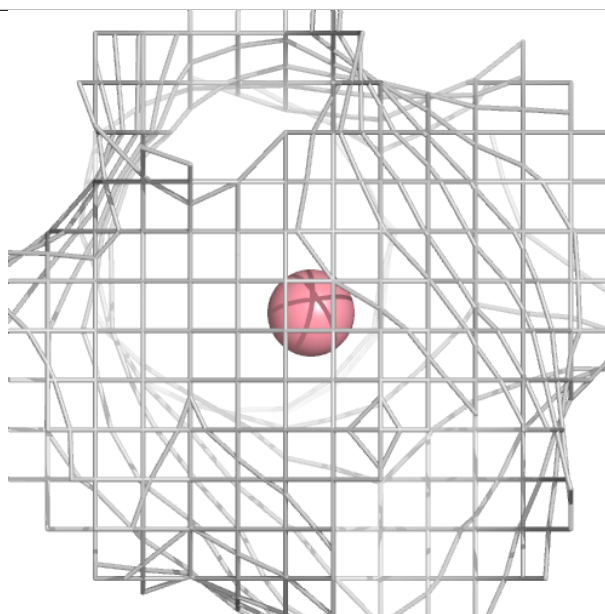
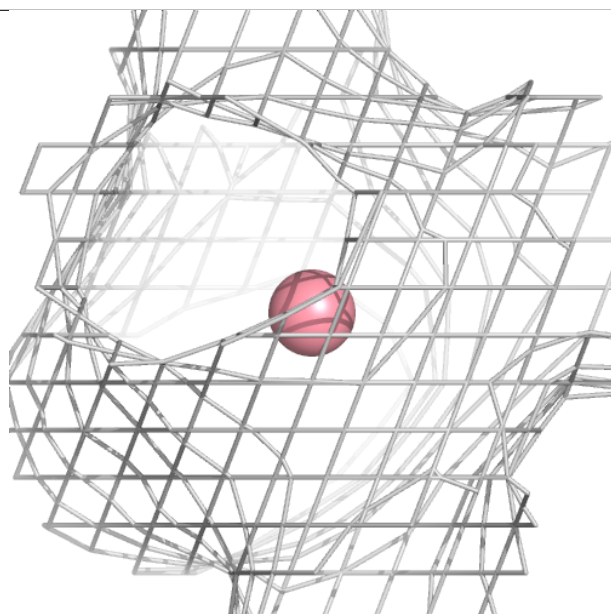
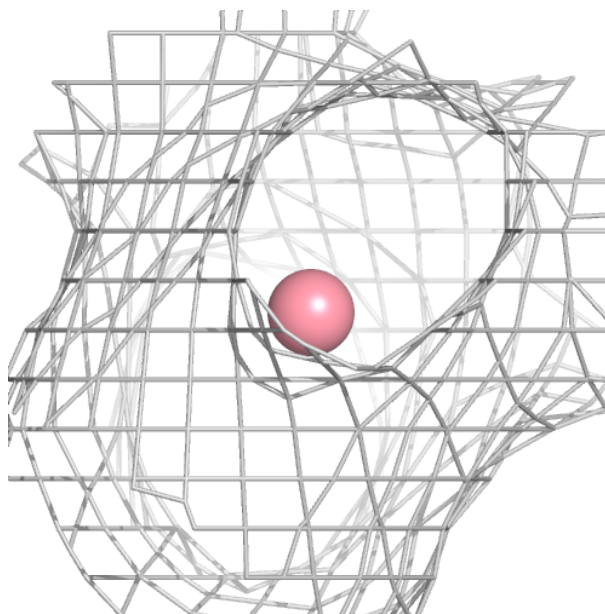
Electron density around CO B 402:

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



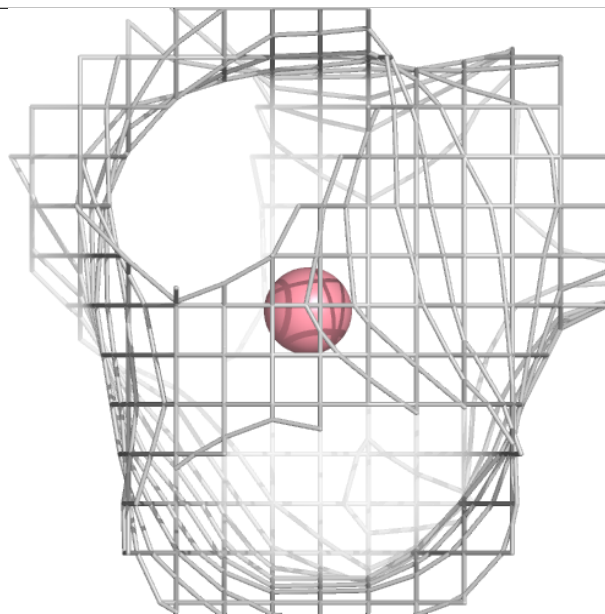
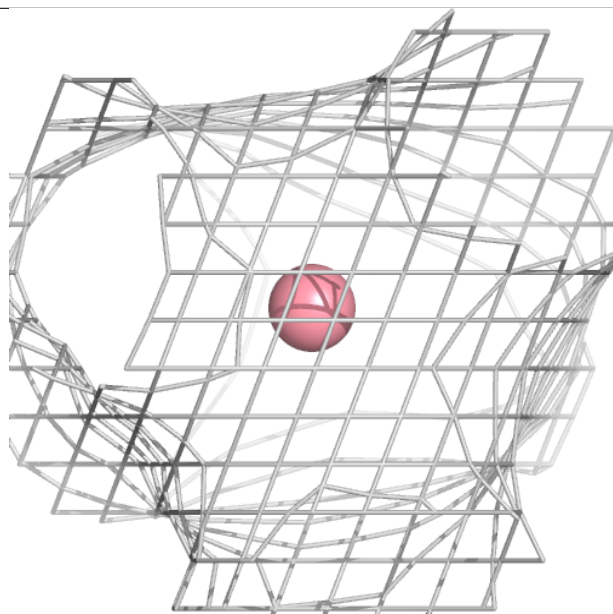
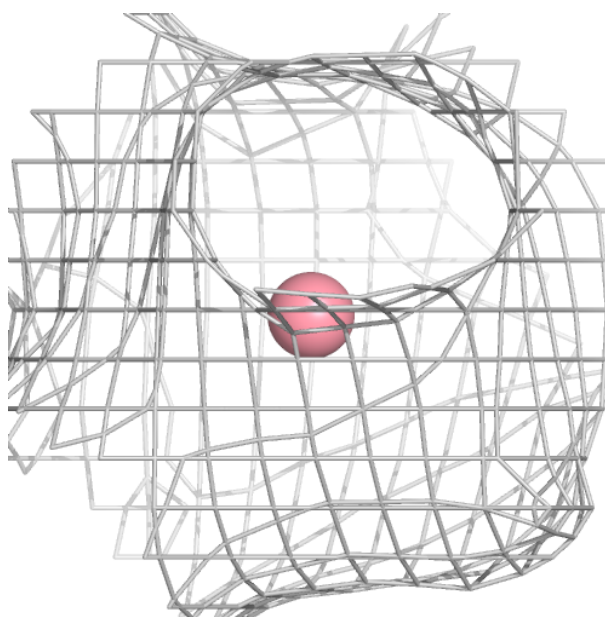
Electron density around CO B 403:

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



Electron density around CO A 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.