



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 3, 2026 – 02:23 AM EDT

PDB ID : 5KOJ / pdb_00005koj
Title : Nitrogenase MoFeP protein in the IDS oxidized state
Authors : Owens, C.P.; Tezcan, F.A.
Deposited on : 2016-06-30
Resolution : 2.59 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (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

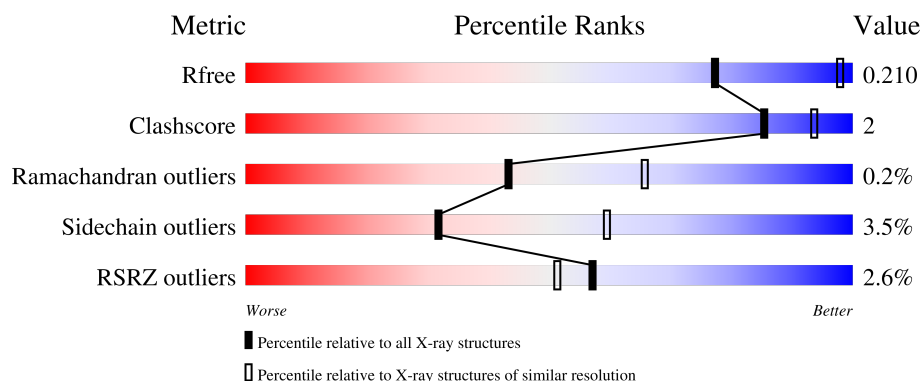
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	499	
1	C	499	
2	B	511	
2	D	511	

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 16070 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.

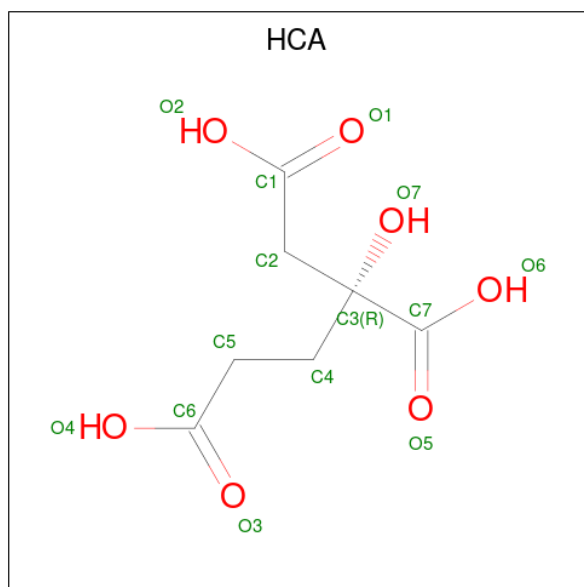
- Molecule 1 is a protein called Nitrogenase protein alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	481	Total	C	N	O	S	0	0	0
			3804	2420	663	701	20			
1	C	480	Total	C	N	O	S	0	0	0
			3800	2418	662	700	20			

- Molecule 2 is a protein called Nitrogenase FeMo beta subunit protein NifK.

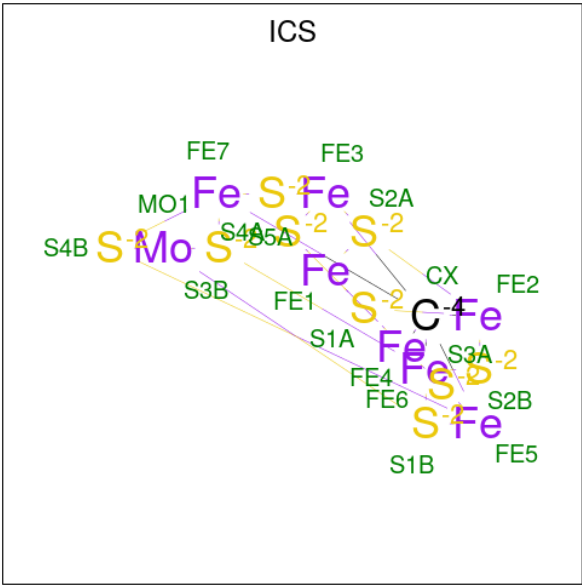
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	510	Total	C	N	O	S	0	0	0
			4006	2547	676	760	23			
2	D	510	Total	C	N	O	S	0	0	0
			4006	2547	676	760	23			

- Molecule 3 is 3-HYDROXY-3-CARBOXY-ADIPIC ACID (CCD ID: HCA) (formula: $C_7H_{10}O_7$).



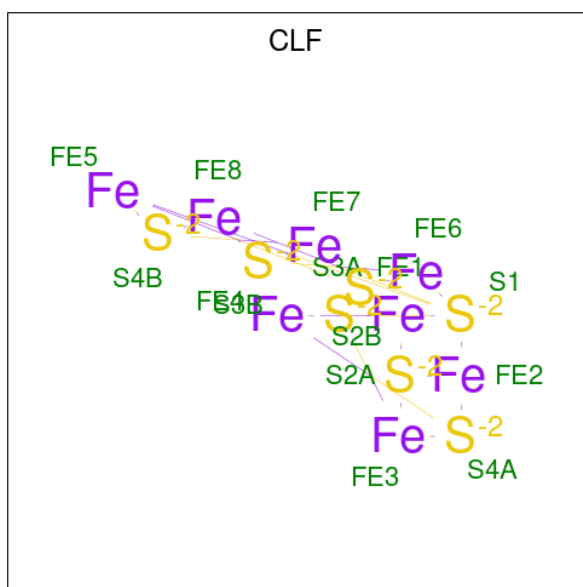
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			14	7	7		
3	C	1	Total	C	O	0	0
			14	7	7		

- Molecule 4 is iron-sulfur-molybdenum cluster with interstitial carbon (CCD ID: ICS) (formula: CFe₇MoS₉).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 18	C 1	Fe 7	Mo 1	S 9	0	0
4	C	1	Total 18	C 1	Fe 7	Mo 1	S 9	0	0

- Molecule 5 is FE(8)-S(7) CLUSTER (CCD ID: CLF) (formula: Fe₈S₇).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Fe S 15 8 7	0	0
5	C	1	Total Fe S 15 8 7	0	0

- Molecule 6 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total Fe 1 1	0	0
6	D	1	Total Fe 1 1	0	0

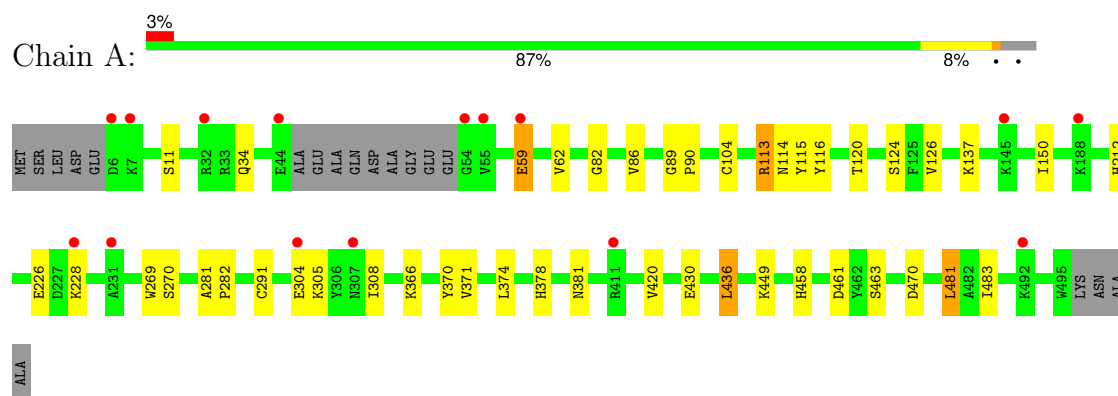
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	87	Total O 87 87	0	0
7	B	93	Total O 93 93	0	0
7	C	76	Total O 76 76	0	0
7	D	102	Total O 102 102	0	0

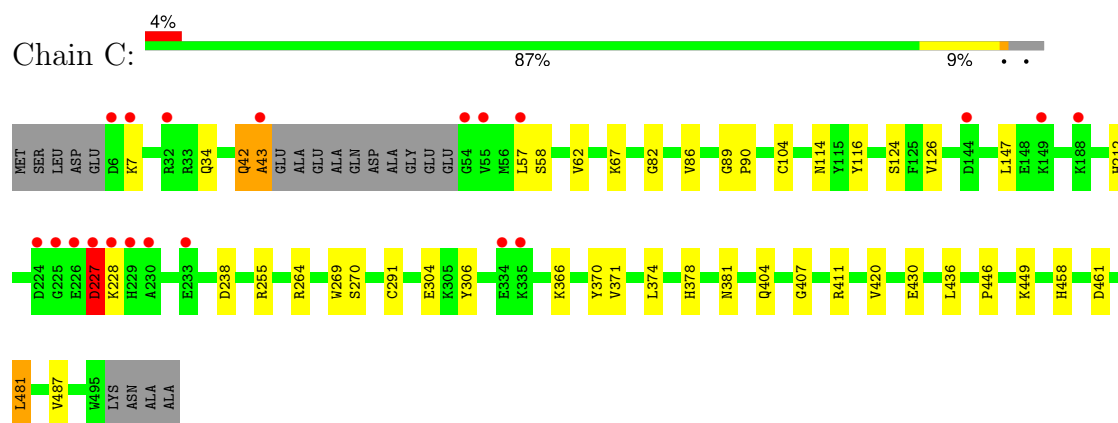
3 Residue-property plots [i](#)

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.

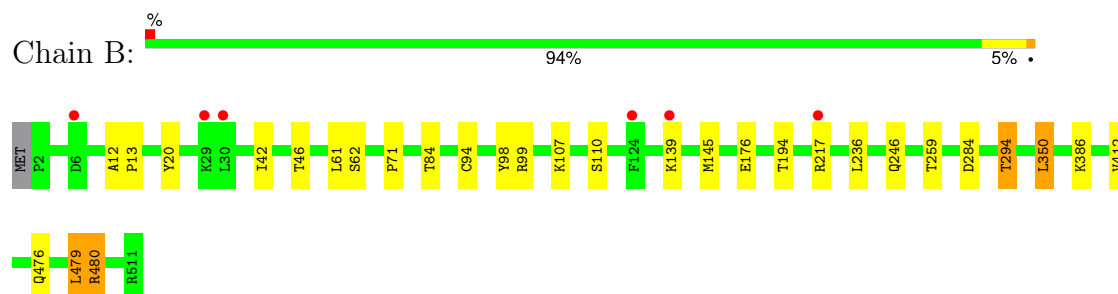
- Molecule 1: Nitrogenase protein alpha chain



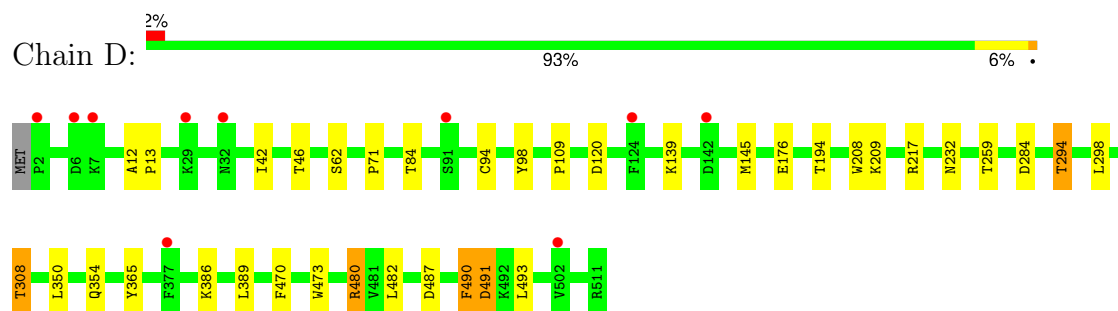
- Molecule 1: Nitrogenase protein alpha chain



- Molecule 2: Nitrogenase FeMo beta subunit protein NifK



● Molecule 2: Nitrogenase FeMo beta subunit protein NifK



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	201.57Å 201.57Å 132.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.59 – 2.59 39.59 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.5 (39.59-2.59) 99.6 (39.59-2.59)	Depositor EDS
R_{merge}	0.41	Depositor
R_{sym}	0.42	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.02 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.168 , 0.208 0.175 , 0.210	Depositor DCC
R_{free} test set	4244 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.335	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 26.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	16070	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.14% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ICS, HCA, CLF, FE

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	0.93	1/3896 (0.0%)	1.02	1/5263 (0.0%)
1	C	0.93	3/3892 (0.1%)	1.05	7/5257 (0.1%)
2	B	0.93	0/4109	0.98	3/5566 (0.1%)
2	D	0.93	1/4109 (0.0%)	1.00	3/5566 (0.1%)
All	All	0.93	5/16006 (0.0%)	1.01	14/21652 (0.1%)

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	C	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	490	PHE	CA-C	-7.09	1.43	1.52
1	C	58	SER	N-CA	6.09	1.54	1.46
1	C	43	ALA	N-CA	5.39	1.56	1.46
1	A	228	LYS	CA-C	5.04	1.59	1.52
1	C	7	LYS	CA-C	5.02	1.58	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	490	PHE	N-CA-C	-9.74	94.81	110.20
1	C	306	TYR	N-CA-C	-9.25	96.08	110.42
1	C	42	GLN	N-CA-C	6.99	122.91	113.97
2	B	412	VAL	N-CA-C	6.80	117.50	110.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	294	THR	CA-C-N	-6.51	112.20	118.97
2	B	294	THR	C-N-CA	-6.51	112.20	118.97
1	C	58	SER	N-CA-C	6.46	124.55	110.80
1	C	306	TYR	CA-C-N	-5.81	113.76	122.24
1	C	306	TYR	C-N-CA	-5.81	113.76	122.24
2	D	294	THR	CA-C-N	-5.47	112.98	119.05
2	D	294	THR	C-N-CA	-5.47	112.98	119.05
1	C	227	ASP	CA-C-N	5.29	131.23	121.70
1	C	227	ASP	C-N-CA	5.29	131.23	121.70
1	A	226	GLU	N-CA-C	-5.08	102.69	110.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	227	ASP	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	3804	0	3710	21	0
1	C	3800	0	3710	19	0
2	B	4006	0	3869	17	0
2	D	4006	0	3869	22	0
3	A	14	0	6	1	0
3	C	14	0	6	1	0
4	A	18	0	0	0	0
4	C	18	0	0	0	0
5	A	15	0	0	0	0
5	C	15	0	0	0	0
6	B	1	0	0	0	0
6	D	1	0	0	0	0
7	A	87	0	0	2	0
7	B	93	0	0	1	0
7	C	76	0	0	2	0
7	D	102	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	16070	0	15170	69	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 2.

All (69) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:490:PHE:O	2:D:491:ASP:CB	2.11	0.98
2:D:490:PHE:O	2:D:491:ASP:HB3	1.79	0.81
1:C:212:HIS:NE2	7:C:601:HOH:O	1.93	0.76
2:D:490:PHE:O	2:D:491:ASP:HB2	1.90	0.72
2:B:99:ARG:HD2	2:B:110:SER:O	1.92	0.69
1:A:212:HIS:NE2	7:A:601:HOH:O	1.93	0.66
2:D:94:CYS:HB3	2:D:98:TYR:CE2	2.34	0.63
2:B:94:CYS:HB3	2:B:98:TYR:CE2	2.35	0.61
1:A:113:ARG:NH2	1:A:463:SER:O	2.35	0.59
1:C:42:GLN:O	1:C:43:ALA:CB	2.52	0.57
1:A:449:LYS:HE3	2:B:259:THR:OG1	2.07	0.55
1:C:227:ASP:CB	1:C:228:LYS:HA	2.36	0.55
1:C:458:HIS:HB3	3:C:501:HCA:O5	2.06	0.55
1:C:449:LYS:HE3	2:D:259:THR:OG1	2.07	0.53
2:D:298:LEU:HD13	2:D:308:THR:HG23	1.90	0.53
2:D:480:ARG:HB3	7:D:788:HOH:O	2.09	0.53
1:C:449:LYS:CE	2:D:259:THR:OG1	2.58	0.52
1:C:481:LEU:C	1:C:481:LEU:HD12	2.35	0.51
1:A:113:ARG:NH1	1:A:115:TYR:OH	2.44	0.51
1:A:449:LYS:CE	2:B:259:THR:OG1	2.59	0.51
2:D:487:ASP:O	2:D:490:PHE:O	2.29	0.51
2:B:480:ARG:HH22	2:D:480:ARG:HH22	1.58	0.50
1:C:116:TYR:CE1	1:C:126:VAL:HB	2.46	0.50
2:D:42:ILE:O	2:D:46:THR:HG23	2.12	0.50
2:B:42:ILE:O	2:B:46:THR:HG23	2.12	0.49
1:A:481:LEU:C	1:A:481:LEU:HD12	2.38	0.49
2:B:107:LYS:HG2	7:B:747:HOH:O	2.11	0.49
1:A:82:GLY:O	1:A:86:VAL:HB	2.12	0.49
1:C:57:LEU:O	1:C:407:GLY:O	2.31	0.48
1:A:458:HIS:HB3	3:A:501:HCA:O5	2.14	0.48
1:C:238:ASP:OD1	1:C:264:ARG:HD3	2.13	0.48
1:C:82:GLY:O	1:C:86:VAL:HB	2.14	0.47
1:A:59:GLU:HA	1:A:59:GLU:OE2	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:350:LEU:HD13	1:C:487:VAL:HG23	1.96	0.47
1:A:116:TYR:CE1	1:A:126:VAL:HB	2.49	0.47
1:A:291:CYS:HA	1:A:374:LEU:HD22	1.97	0.46
1:C:291:CYS:HA	1:C:374:LEU:HD22	1.96	0.46
2:B:476:GLN:CD	2:B:480:ARG:HH21	2.23	0.45
2:D:354:GLN:HB2	7:D:757:HOH:O	2.15	0.45
2:B:217:ARG:NH1	2:B:284:ASP:OD2	2.49	0.44
1:C:255:ARG:NH2	7:C:606:HOH:O	2.51	0.44
1:C:446:PRO:HG3	2:D:109:PRO:HB3	1.99	0.44
1:A:116:TYR:OH	1:A:120:THR:OG1	2.27	0.44
2:B:94:CYS:HB3	2:B:98:TYR:CZ	2.53	0.44
1:A:269:TRP:HA	1:A:270:SER:HA	1.78	0.44
2:D:71:PRO:HB2	2:D:98:TYR:CZ	2.53	0.44
1:A:150:ILE:HG23	2:B:61:LEU:HD13	2.00	0.43
1:C:89:GLY:N	1:C:90:PRO:CD	2.82	0.43
1:A:89:GLY:N	1:A:90:PRO:CD	2.82	0.43
2:B:12:ALA:HB3	2:B:13:PRO:CD	2.49	0.43
2:D:94:CYS:HB3	2:D:98:TYR:CZ	2.53	0.43
2:B:479:LEU:HD13	2:D:470:PHE:CZ	2.54	0.43
2:D:12:ALA:HB3	2:D:13:PRO:CD	2.49	0.43
1:A:212:HIS:CE1	7:A:601:HOH:O	2.60	0.42
2:D:208:TRP:O	2:D:209:LYS:C	2.63	0.42
1:A:370:TYR:CZ	1:A:420:VAL:HG12	2.55	0.42
2:B:71:PRO:HB2	2:B:98:TYR:CZ	2.55	0.42
1:C:370:TYR:CZ	1:C:420:VAL:HG12	2.55	0.42
1:A:436:LEU:HG	1:A:483:ILE:HD11	2.01	0.42
2:D:217:ARG:NH1	2:D:284:ASP:OD2	2.50	0.41
1:A:281:ALA:N	1:A:282:PRO:CD	2.82	0.41
2:B:84:THR:OG1	2:B:145:MET:HE2	2.21	0.41
1:C:269:TRP:HA	1:C:270:SER:HA	1.77	0.41
2:D:365:TYR:HA	2:D:389:LEU:O	2.21	0.41
1:A:470:ASP:OD2	2:B:20:TYR:OH	2.34	0.40
1:C:42:GLN:O	1:C:43:ALA:HB2	2.21	0.40
1:A:282:PRO:O	1:A:308:ILE:HD11	2.22	0.40
2:D:84:THR:OG1	2:D:145:MET:HE2	2.21	0.40
2:D:232:ASN:HB3	2:D:473:TRP:CE3	2.57	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	477/499 (96%)	455 (95%)	22 (5%)	0	100	100
1	C	476/499 (95%)	453 (95%)	22 (5%)	1 (0%)	43	66
2	B	508/511 (99%)	496 (98%)	11 (2%)	1 (0%)	43	66
2	D	508/511 (99%)	497 (98%)	9 (2%)	2 (0%)	30	51
All	All	1969/2020 (98%)	1901 (96%)	64 (3%)	4 (0%)	43	66

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	227	ASP
2	D	491	ASP
2	B	294	THR
2	D	294	THR

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	398/413 (96%)	379 (95%)	19 (5%)	23	47
1	C	398/413 (96%)	380 (96%)	18 (4%)	24	50
2	B	427/428 (100%)	417 (98%)	10 (2%)	44	71
2	D	427/428 (100%)	416 (97%)	11 (3%)	40	68
All	All	1650/1682 (98%)	1592 (96%)	58 (4%)	32	59

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	SER
1	A	34	GLN
1	A	59	GLU
1	A	62	VAL
1	A	104	CYS
1	A	113	ARG
1	A	114	ASN
1	A	124	SER
1	A	137	LYS
1	A	304	GLU
1	A	305	LYS
1	A	366	LYS
1	A	371	VAL
1	A	378	HIS
1	A	381	ASN
1	A	430	GLU
1	A	436	LEU
1	A	461	ASP
1	A	481	LEU
2	B	62	SER
2	B	139	LYS
2	B	176	GLU
2	B	194	THR
2	B	236	LEU
2	B	246	GLN
2	B	350	LEU
2	B	386	LYS
2	B	479	LEU
2	B	480	ARG
1	C	34	GLN
1	C	62	VAL
1	C	67	LYS
1	C	104	CYS
1	C	114	ASN
1	C	124	SER
1	C	147	LEU
1	C	304	GLU
1	C	366	LYS
1	C	371	VAL
1	C	378	HIS
1	C	381	ASN
1	C	404	GLN

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Mol	Chain	Res	Type
1	C	411	ARG
1	C	430	GLU
1	C	436	LEU
1	C	461	ASP
1	C	481	LEU
2	D	62	SER
2	D	120	ASP
2	D	139	LYS
2	D	176	GLU
2	D	194	THR
2	D	308	THR
2	D	350	LEU
2	D	386	LYS
2	D	480	ARG
2	D	482	LEU
2	D	493	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	400	ASN
2	B	21	GLN
2	B	358	HIS
1	C	129	GLN
1	C	378	HIS
1	C	448	GLN
2	D	358	HIS
2	D	476	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 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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	HCA	A	501	4	13,13,13	0.86	0	15,18,18	1.86	4 (26%)
4	ICS	A	502	1,3	6,30,30	1.29	0	-		
3	HCA	C	501	4	13,13,13	0.97	0	15,18,18	1.45	4 (26%)
5	CLF	A	503	1,2	0,24,24	-	-	-		
5	CLF	C	503	1,2	0,24,24	-	-	-		
4	ICS	C	502	1,3	6,30,30	1.93	3 (50%)	-		

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	HCA	A	501	4	-	3/17/17/17	-
3	HCA	C	501	4	-	6/17/17/17	-
5	CLF	C	503	1,2	-	-	0/12/10/10
5	CLF	A	503	1,2	-	-	0/12/10/10

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	502	ICS	S3A-FE5	3.07	2.32	2.24
4	C	502	ICS	S5A-FE3	-2.14	2.19	2.24
4	C	502	ICS	S2B-FE6	-2.11	2.19	2.24

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	HCA	O6-C7-C3	5.02	122.78	113.14
3	A	501	HCA	O3-C6-C5	-3.12	113.19	123.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	501	HCA	O7-C3-C7	-2.70	105.13	108.96
3	A	501	HCA	O4-C6-C5	2.34	121.38	114.00
3	C	501	HCA	O1-C1-C2	-2.33	116.36	122.95
3	A	501	HCA	O5-C7-C3	-2.23	117.77	122.09
3	C	501	HCA	O4-C6-C5	2.06	120.52	114.00
3	C	501	HCA	O2-C1-C2	2.06	120.86	114.35

There are no chirality outliers.

All (9) torsion outliers are listed below:

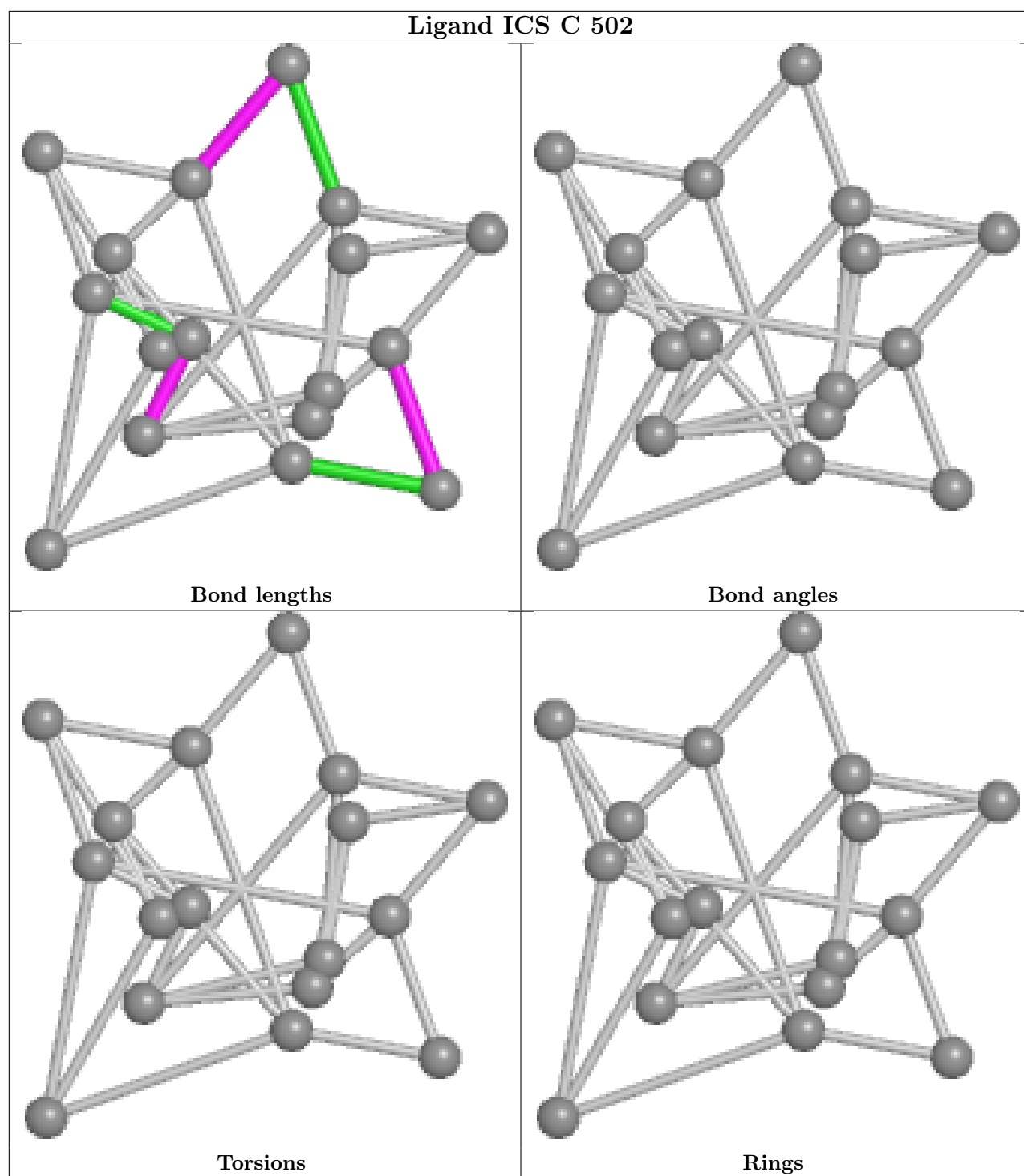
Mol	Chain	Res	Type	Atoms
3	A	501	HCA	C2-C3-C4-C5
3	A	501	HCA	C7-C3-C4-C5
3	A	501	HCA	O7-C3-C4-C5
3	C	501	HCA	C2-C3-C4-C5
3	C	501	HCA	C7-C3-C4-C5
3	C	501	HCA	O7-C3-C4-C5
3	C	501	HCA	C3-C4-C5-C6
3	C	501	HCA	C4-C5-C6-O4
3	C	501	HCA	C4-C5-C6-O3

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	HCA	1	0
3	C	501	HCA	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 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 ⓘ

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	481/499 (96%)	0.43	15 (3%) 51 45	28, 38, 61, 79	0
1	C	480/499 (96%)	0.53	20 (4%) 40 35	28, 39, 64, 100	0
2	B	510/511 (99%)	0.23	6 (1%) 76 73	25, 36, 56, 76	0
2	D	510/511 (99%)	0.31	10 (1%) 65 60	25, 37, 57, 86	0
All	All	1981/2020 (98%)	0.37	51 (2%) 57 51	25, 38, 59, 100	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	44	GLU	5.1
1	A	54	GLY	4.7
2	D	124	PHE	4.3
1	C	6	ASP	4.2
1	C	54	GLY	4.0
1	C	227	ASP	3.8
1	C	43	ALA	3.8
1	C	225	GLY	3.8
1	A	32	ARG	3.3
1	A	59	GLU	3.0
1	C	188	LYS	3.0
2	D	91	SER	2.9
1	A	55	VAL	2.8
1	A	492	LYS	2.7
1	A	228	LYS	2.6
1	C	228	LYS	2.6
2	B	139	LYS	2.6
1	C	32	ARG	2.5
1	C	230	ALA	2.5
2	D	142	ASP	2.5
2	D	377	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	217	ARG	2.5
2	B	6	ASP	2.4
2	B	30	LEU	2.4
1	C	7	LYS	2.4
1	C	149	LYS	2.4
2	B	124	PHE	2.4
2	D	502	VAL	2.4
1	C	226	GLU	2.4
1	A	7	LYS	2.4
1	C	57	LEU	2.3
1	A	411	ARG	2.3
2	B	29	LYS	2.3
2	D	7	LYS	2.3
1	A	231	ALA	2.3
1	C	229	HIS	2.2
1	C	224	ASP	2.2
1	C	233	GLU	2.2
1	C	55	VAL	2.2
2	D	2	PRO	2.2
1	A	307	ASN	2.2
2	D	32	ASN	2.2
1	C	334	GLU	2.2
2	D	6	ASP	2.2
1	A	188	LYS	2.1
1	A	6	ASP	2.1
1	A	145	LYS	2.1
1	C	144	ASP	2.1
1	C	335	LYS	2.1
1	A	304	GLU	2.0
2	D	29	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands

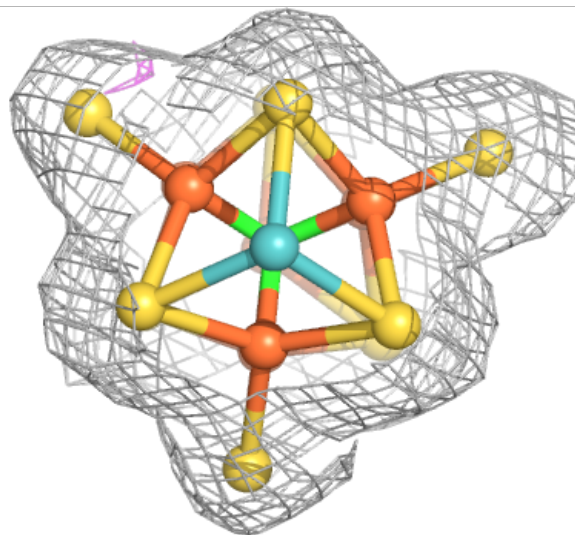
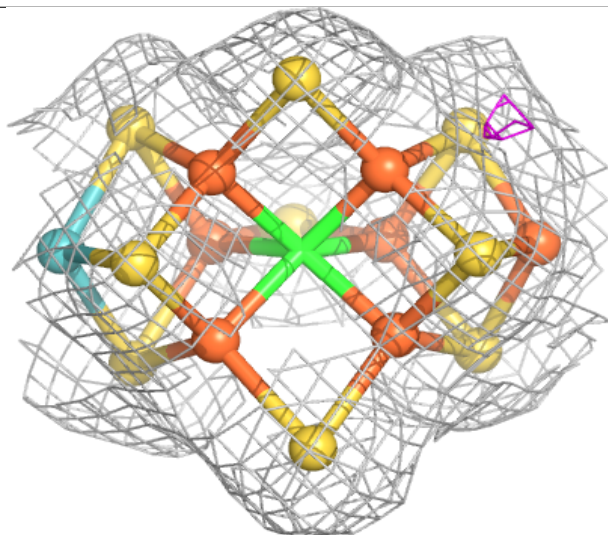
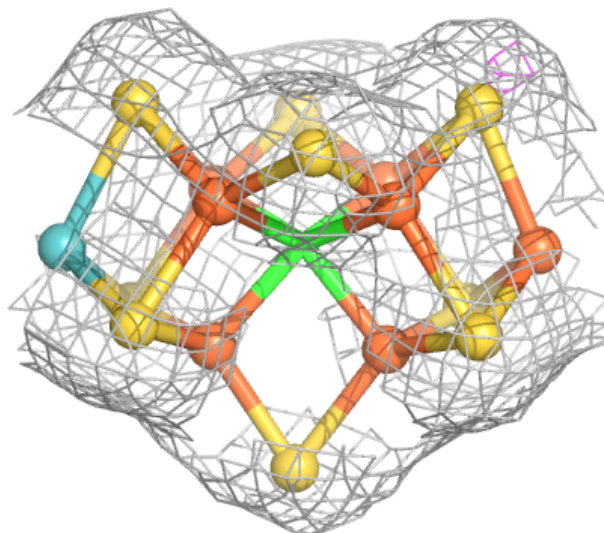
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	FE	B	601	1/1	0.89	0.11	81,81,81,81	0
3	HCA	A	501	14/14	0.95	0.09	28,33,38,43	0
5	CLF	C	503	15/15	0.97	0.05	34,38,49,54	0
3	HCA	C	501	14/14	0.97	0.08	31,37,39,43	0
6	FE	D	601	1/1	0.97	0.06	62,62,62,62	0
5	CLF	A	503	15/15	0.98	0.04	28,34,38,44	2
4	ICS	A	502	18/18	0.99	0.03	27,32,34,37	0
4	ICS	C	502	18/18	0.99	0.03	31,34,38,40	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 ICS C 502:

$2mF_o-DF_c$ (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.