



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 8, 2022 – 03:29 PM EST

PDB ID : 8E3U  
Title : Nickel-reconstituted nitrogenase MoFeP mutant S188A from *Azotobacter vinelandii* after IDS oxidation  
Authors : Rutledge, H.L.; Tezcan, F.A.  
Deposited on : 2022-08-17  
Resolution : 1.99 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.2

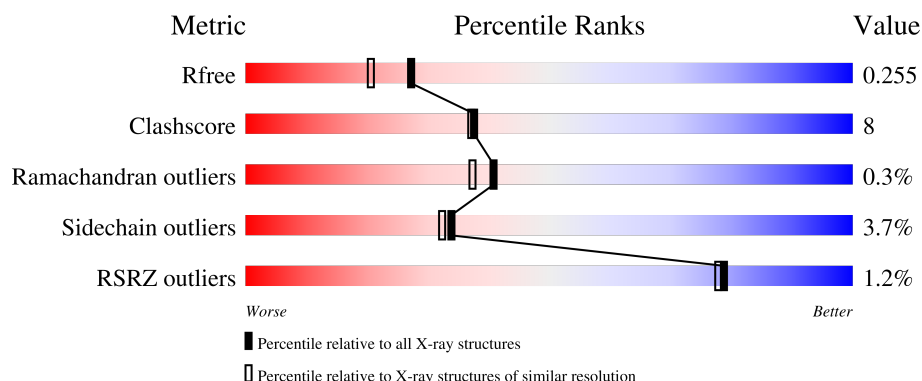
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.99 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	492	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>18%</div> <div>• •</div> </div> </div>
1	C	492	<div> <div>2%</div> <div> <div></div> <div>77%</div> <div>16%</div> <div>• 5%</div> </div> </div>
2	B	523	<div> <div></div> <div> <div></div> <div>84%</div> <div>15%</div> <div>•</div> </div> </div>
2	D	523	<div> <div></div> <div> <div></div> <div>82%</div> <div>17%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	UFF	A	601	-	-	X	-

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 31928 atoms, of which 15166 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 molybdenum-iron protein alpha chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	473	Total	C	H	N	O	S	0	0	0
			7277	2359	3570	628	696	24			
1	C	469	Total	C	H	N	O	S	0	0	0
			7219	2341	3546	625	683	24			

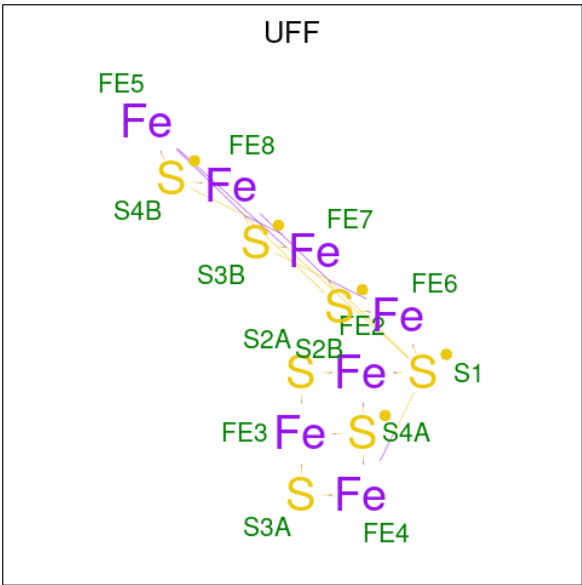
- Molecule 2 is a protein called Nitrogenase molybdenum-iron protein beta chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	522	Total	C	H	N	O	S	0	0	0
			8167	2647	4025	699	768	28			
2	D	522	Total	C	H	N	O	S	0	0	0
			8148	2644	4013	697	766	28			

There are 2 discrepancies between the modelled and reference sequences:

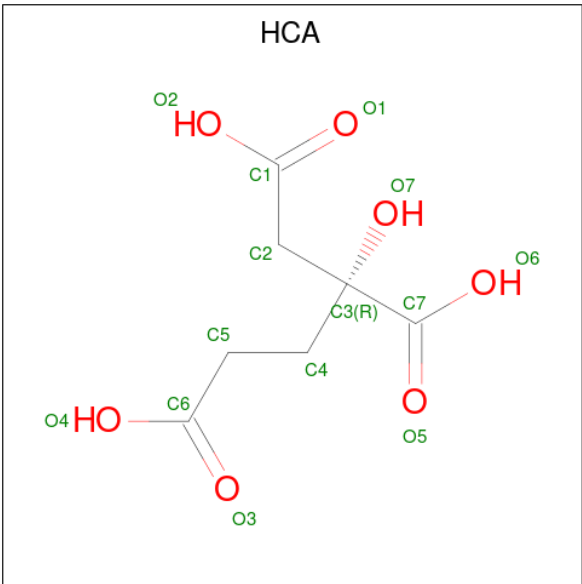
Chain	Residue	Modelled	Actual	Comment	Reference
B	188	ALA	SER	engineered mutation	UNP C1DGZ8
D	188	ALA	SER	engineered mutation	UNP C1DGZ8

- Molecule 3 is FE(7)-S(7) CLUSTER (three-letter code: UFF) (formula: Fe<sub>7</sub>S<sub>7</sub>) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is 3-HYDROXY-3-CARBOXY-ADIPIC ACID (three-letter code: HCA) (formula: C<sub>7</sub>H<sub>10</sub>O<sub>7</sub>).



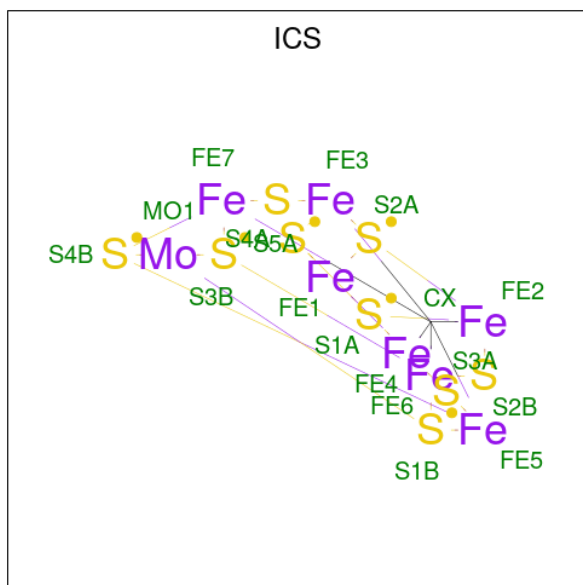
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			20	7	6	7		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	H	O	0	0
			20	7	6	7		

- Molecule 5 is iron-sulfur-molybdenum cluster with interstitial carbon (three-letter code: ICS) (formula:  $\text{CFe}_7\text{MoS}_9$ ).



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

- Molecule 6 is FE (III) ION (three-letter code: 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 SODIUM ION (three-letter code: NA) (formula: Na).

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

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	D	1	Total 1	Na 1	0	0

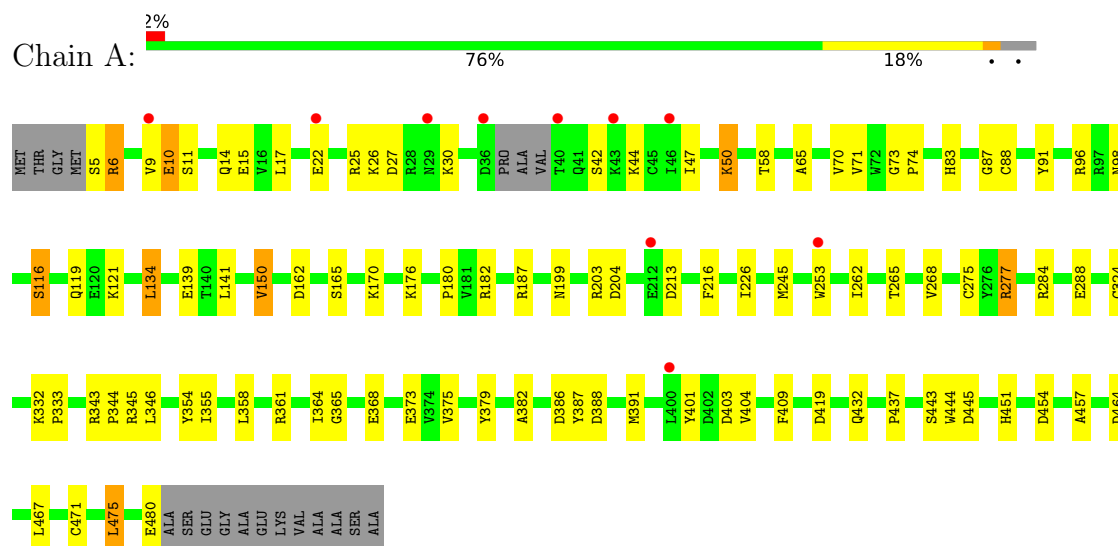
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	208	Total 208	O 208	0	0
8	B	317	Total 317	O 317	0	0
8	C	200	Total 200	O 200	0	0
8	D	284	Total 284	O 284	0	0

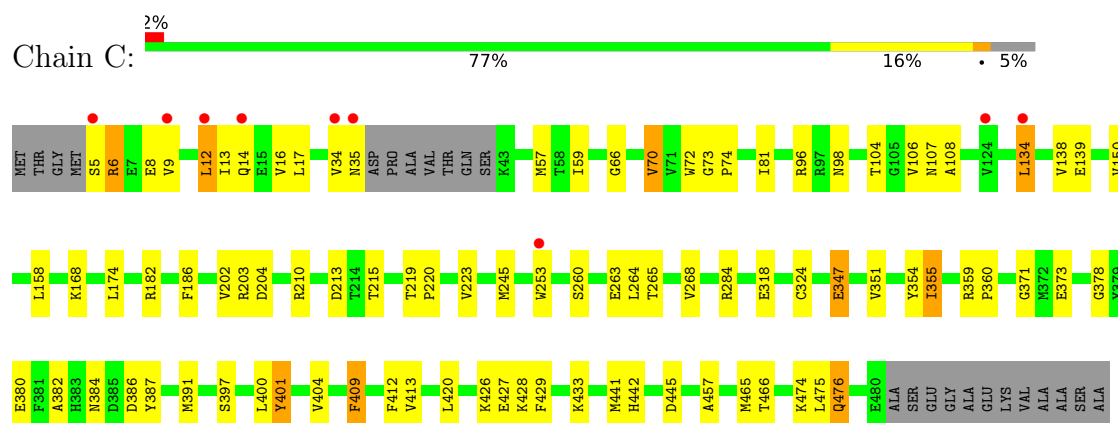
### 3 Residue-property plots [i](#)

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.

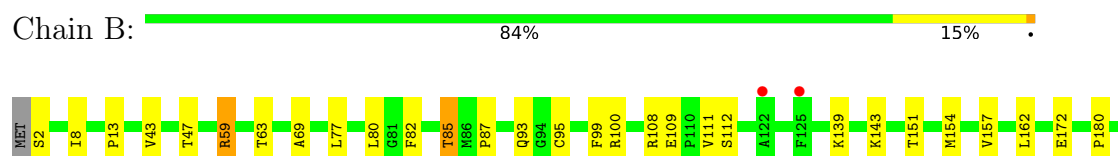
- Molecule 1: Nitrogenase molybdenum-iron protein alpha chain



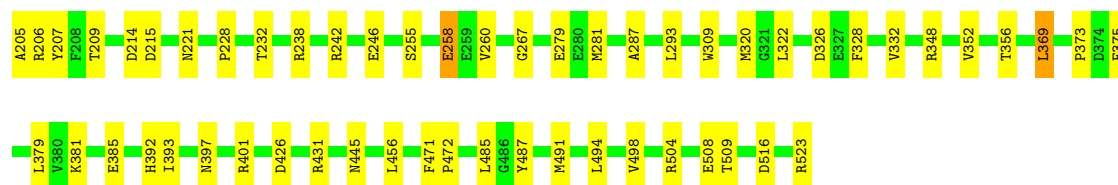
- Molecule 1: Nitrogenase molybdenum-iron protein alpha chain



- Molecule 2: Nitrogenase molybdenum-iron protein beta chain

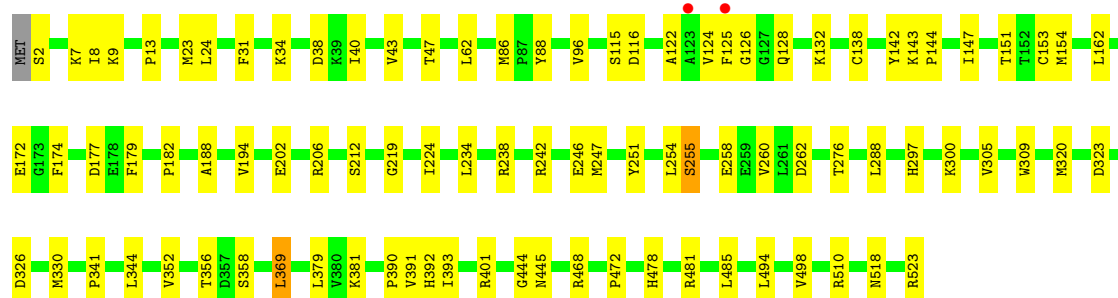






• Molecule 2: Nitrogenase molybdenum-iron protein beta chain

Chain D: 82% 17%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.26Å 129.01Å 107.76Å 90.00° 109.22° 90.00°	Depositor
Resolution (Å)	79.89 – 1.99 79.89 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.0 (79.89-1.99) 99.6 (79.89-1.99)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.18 (at 1.98Å)	Xtriage
Refinement program	PHENIX 1.19.1-4122	Depositor
R, $R_{free}$	0.202 , 0.253 0.205 , 0.255	Depositor DCC
$R_{free}$ test set	6860 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.2	Xtriage
Anisotropy	0.450	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 41.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.025 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	31928	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.53% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.82	0/3793	0.87	0/5124
1	C	0.80	0/3759	0.88	0/5078
2	B	0.87	0/4248	0.90	2/5751 (0.0%)
2	D	0.87	0/4241	0.91	0/5743
All	All	0.84	0/16041	0.89	2/21696 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	59	ARG	NE-CZ-NH1	5.29	122.94	120.30
2	B	523	ARG	NE-CZ-NH1	5.18	122.89	120.30

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	3707	3570	3585	69	0
1	C	3673	3546	3558	62	0
2	B	4142	4025	4020	56	0
2	D	4135	4013	4004	63	0
3	A	14	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	14	0	0	0	0
4	A	14	6	6	0	0
4	C	14	6	6	0	0
5	A	18	0	0	1	0
5	C	18	0	0	2	0
6	B	1	0	0	0	0
6	D	1	0	0	0	0
7	B	1	0	0	1	0
7	D	1	0	0	0	0
8	A	208	0	0	8	0
8	B	317	0	0	6	0
8	C	200	0	0	2	0
8	D	284	0	0	4	0
All	All	16762	15166	15179	237	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 8.

All (237) 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:B:209:THR:HG21	2:B:309:TRP:HE1	1.26	0.96
2:B:209:THR:HG21	2:B:309:TRP:NE1	1.89	0.86
3:A:601:UFF:FE5	3:A:601:UFF:S2B	1.71	0.83
2:B:143:LYS:HE2	2:B:143:LYS:HA	1.59	0.82
1:C:134:LEU:HD13	2:D:62:LEU:HD13	1.68	0.76
1:A:457:ALA:HB1	2:B:8:ILE:HD12	1.68	0.75
1:C:457:ALA:HB1	2:D:8:ILE:HD12	1.68	0.74
1:C:35:ASN:HB2	1:C:400:LEU:HD11	1.69	0.72
2:D:518:ASN:O	2:D:523:ARG:NH2	2.24	0.71
2:B:348:ARG:NH2	1:C:474:LYS:O	2.24	0.71
1:C:12:LEU:HD23	1:C:13:ILE:N	2.07	0.69
3:A:601:UFF:S3A	7:B:602:NA:NA	2.08	0.67
2:B:494:LEU:O	2:B:498:VAL:HG23	1.95	0.66
2:D:238:ARG:HG3	2:D:238:ARG:HH11	1.61	0.66
2:B:205:ALA:O	2:B:209:THR:HB	1.97	0.65
2:B:238:ARG:NH1	2:B:258:GLU:OE1	2.32	0.62
1:A:253:TRP:CZ2	1:A:262:ILE:HG23	2.35	0.62
1:C:354:TYR:CZ	1:C:404:VAL:HG12	2.35	0.62
1:A:284:ARG:HG2	1:A:284:ARG:HH11	1.65	0.61
1:A:354:TYR:CZ	1:A:404:VAL:HG12	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:364:ILE:O	1:A:368:GLU:HG3	2.02	0.59
1:A:277:ARG:HD3	8:A:725:HOH:O	2.03	0.59
1:C:203:ARG:HD2	1:C:204:ASP:OD1	2.02	0.59
2:D:254:LEU:HA	2:D:276:THR:HG21	1.85	0.59
1:C:96:ARG:NH1	5:C:603:ICS:S5A	2.77	0.58
1:A:6:ARG:H	1:A:9:VAL:HG23	1.68	0.58
1:A:432:GLN:NE2	8:A:704:HOH:O	2.30	0.58
1:C:476:GLN:HE21	1:C:476:GLN:HA	1.69	0.57
1:A:11:SER:O	1:A:15:GLU:OE1	2.22	0.57
2:D:124:VAL:HG12	2:D:125:PHE:CD1	2.40	0.57
2:D:124:VAL:HG12	2:D:125:PHE:CE1	2.40	0.57
1:C:284:ARG:HH11	1:C:284:ARG:HG2	1.70	0.57
1:A:134:LEU:C	1:A:134:LEU:HD12	2.26	0.56
1:C:426:LYS:NZ	8:C:708:HOH:O	2.38	0.56
1:A:365:GLY:N	8:A:711:HOH:O	2.38	0.56
2:D:238:ARG:HH11	2:D:238:ARG:CG	2.19	0.55
1:C:409:PHE:O	1:C:413:VAL:HG23	2.07	0.55
1:A:265:THR:O	1:A:268:VAL:HG22	2.06	0.55
2:B:139:LYS:O	2:B:143:LYS:HE3	2.05	0.55
1:A:203:ARG:HD2	1:A:204:ASP:OD1	2.07	0.55
2:B:77:LEU:HA	2:B:80:LEU:HD12	1.89	0.54
1:C:35:ASN:HB2	1:C:400:LEU:CD1	2.37	0.54
2:B:242:ARG:HD3	2:B:246:GLU:OE2	2.08	0.54
1:C:81:ILE:HG21	1:C:134:LEU:HD21	1.90	0.54
1:A:42:SER:C	1:A:44:LYS:H	2.09	0.54
2:D:494:LEU:C	2:D:494:LEU:HD23	2.29	0.54
2:D:122:ALA:O	2:D:126:GLY:N	2.40	0.54
1:C:351:VAL:HA	1:C:420:LEU:O	2.08	0.53
1:A:47:ILE:HD12	1:A:50:LYS:CD	2.38	0.53
1:A:88:CYS:SG	3:A:601:UFF:S2B	3.06	0.53
1:A:17:LEU:HB3	1:A:25:ARG:HG3	1.90	0.53
2:D:369:LEU:N	2:D:369:LEU:CD1	2.71	0.53
2:D:369:LEU:N	2:D:369:LEU:HD13	2.23	0.53
1:C:265:THR:O	1:C:268:VAL:HG22	2.09	0.53
2:D:320:MET:HG3	2:D:485:LEU:HD23	1.90	0.53
2:B:352:VAL:O	2:B:356:THR:HG23	2.07	0.53
2:D:352:VAL:O	2:D:356:THR:HG23	2.07	0.53
1:C:9:VAL:O	1:C:13:ILE:HD12	2.08	0.53
1:C:12:LEU:O	1:C:16:VAL:HG23	2.08	0.53
1:C:476:GLN:HA	1:C:476:GLN:NE2	2.24	0.53
1:C:134:LEU:HD12	1:C:134:LEU:O	2.07	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:320:MET:HG3	2:B:485:LEU:HD23	1.89	0.52
2:B:206:ARG:HD2	8:B:832:HOH:O	2.09	0.52
2:D:391:VAL:HG12	2:D:392:HIS:CE1	2.45	0.52
1:C:34:VAL:HG12	1:C:397:SER:HA	1.92	0.52
1:C:134:LEU:HD12	1:C:134:LEU:C	2.31	0.52
1:A:382:ALA:HB1	1:A:386:ASP:HB2	1.92	0.51
2:B:369:LEU:HD21	2:B:393:ILE:HG23	1.93	0.51
2:D:510:ARG:HD3	8:D:775:HOH:O	2.10	0.51
1:A:284:ARG:HG2	1:A:284:ARG:NH1	2.26	0.51
2:B:180:PRO:HA	2:B:207:TYR:OH	2.11	0.51
2:B:100:ARG:HD2	2:B:111:VAL:O	2.11	0.51
2:D:369:LEU:HD13	2:D:369:LEU:H	1.76	0.50
2:B:151:THR:HG23	2:B:162:LEU:HD11	1.92	0.50
2:D:147:ILE:HG13	2:D:179:PHE:HE2	1.77	0.50
2:D:358:SER:HB3	2:D:498:VAL:HG21	1.92	0.50
2:B:456:LEU:HD23	2:D:510:ARG:HG2	1.94	0.50
2:B:369:LEU:HD13	2:B:392:HIS:O	2.11	0.50
1:C:158:LEU:HG	1:C:186:PHE:CE2	2.47	0.50
1:C:245:MET:HG3	1:C:324:CYS:HA	1.94	0.49
1:A:119:GLN:HB2	1:A:121:LYS:HE3	1.93	0.49
2:B:232:THR:HG21	2:B:471:PHE:CD1	2.47	0.49
2:D:247:MET:HG2	2:D:341:PRO:HD3	1.94	0.49
2:B:369:LEU:HD22	2:B:369:LEU:O	2.12	0.49
1:A:150:VAL:HG13	1:A:180:PRO:HA	1.93	0.49
2:D:23:MET:HE2	2:D:24:LEU:HG	1.95	0.49
2:D:219:GLY:HA2	2:D:288:LEU:HD23	1.93	0.49
2:B:43:VAL:O	2:B:47:THR:HG23	2.12	0.49
2:D:445:ASN:HB2	2:D:472:PRO:O	2.13	0.49
1:A:5:SER:O	1:A:6:ARG:CB	2.62	0.48
2:B:445:ASN:HB2	2:B:472:PRO:O	2.14	0.48
1:A:332:LYS:N	1:A:333:PRO:HD2	2.28	0.48
1:C:213:ASP:OD1	1:C:215:THR:HG23	2.12	0.48
2:B:369:LEU:HD11	2:B:393:ILE:HG12	1.96	0.48
2:D:444:GLY:O	2:D:468:ARG:HA	2.14	0.47
1:A:47:ILE:HD12	1:A:50:LYS:HD2	1.97	0.47
1:A:91:TYR:OH	2:B:69:ALA:HB1	2.14	0.47
2:B:504:ARG:O	2:B:508:GLU:HG3	2.13	0.47
1:C:12:LEU:HD21	1:C:412:PHE:CD1	2.48	0.47
2:D:43:VAL:O	2:D:47:THR:HG23	2.14	0.47
1:A:162:ASP:OD2	1:A:165:SER:CB	2.62	0.47
2:D:305:VAL:O	2:D:309:TRP:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:343:ARG:N	1:A:344:PRO:CD	2.78	0.47
2:B:221:ASN:OD1	2:B:287:ALA:HA	2.14	0.47
1:C:12:LEU:HD23	1:C:12:LEU:C	2.34	0.47
1:C:66:GLY:O	1:C:70:VAL:HG22	2.15	0.47
1:A:10:GLU:OE1	1:A:10:GLU:HA	2.15	0.47
1:C:134:LEU:CD1	2:D:62:LEU:HD13	2.42	0.47
2:B:375:PHE:CZ	2:B:379:LEU:CD1	2.98	0.47
1:C:12:LEU:HD21	1:C:412:PHE:CE1	2.49	0.47
1:C:355:ILE:HB	1:C:360:PRO:HD3	1.96	0.47
1:A:42:SER:C	1:A:44:LYS:N	2.69	0.47
2:B:63:THR:HG23	2:B:426:ASP:OD2	2.15	0.47
1:C:433:LYS:HD3	8:D:830:HOH:O	2.15	0.46
2:D:9:LYS:HD2	2:D:13:PRO:HG2	1.97	0.46
1:A:216:PHE:N	8:A:721:HOH:O	2.46	0.46
1:C:219:THR:HB	1:C:220:PRO:HD2	1.97	0.46
2:D:147:ILE:O	2:D:182:PRO:HD2	2.16	0.46
1:A:47:ILE:HD12	1:A:50:LYS:HD3	1.97	0.46
2:D:379:LEU:HD23	2:D:379:LEU:HA	1.80	0.46
2:B:157:VAL:O	8:B:701:HOH:O	2.21	0.46
1:A:22:GLU:OE2	1:A:25:ARG:NH1	2.49	0.46
2:B:375:PHE:CZ	2:B:379:LEU:HD11	2.51	0.46
1:C:134:LEU:HD13	2:D:62:LEU:CD1	2.43	0.46
1:C:354:TYR:CE2	1:C:404:VAL:HG12	2.50	0.46
2:B:95:CYS:HB3	2:B:99:PHE:CZ	2.51	0.46
2:D:234:LEU:HD13	2:D:262:ASP:HB2	1.98	0.46
1:A:6:ARG:N	1:A:9:VAL:HG23	2.30	0.45
1:A:346:LEU:HD21	1:A:464:ASP:HA	1.97	0.45
1:C:134:LEU:HD12	1:C:138:VAL:HG23	1.98	0.45
1:C:355:ILE:O	1:C:380:GLU:HG3	2.15	0.45
1:A:42:SER:HB3	1:A:388:ASP:OD1	2.15	0.45
2:B:397:ASN:HB2	8:B:818:HOH:O	2.17	0.45
2:D:128:GLN:O	2:D:132:LYS:HG3	2.17	0.45
1:A:253:TRP:HE1	1:A:265:THR:HG1	1.63	0.45
2:B:82:PHE:O	2:B:85:THR:HG23	2.17	0.45
1:C:59:ILE:HD13	1:C:354:TYR:CE2	2.51	0.45
1:C:210:ARG:HD3	1:C:263:GLU:HB3	1.98	0.45
1:C:359:ARG:N	1:C:360:PRO:CD	2.78	0.45
2:B:328:PHE:O	2:B:332:VAL:HG23	2.15	0.45
1:A:87:GLY:HA3	3:A:601:UFF:S2B	2.57	0.45
1:A:170:LYS:HD2	1:A:170:LYS:HA	1.81	0.45
1:A:141:LEU:CD1	2:B:59:ARG:HD2	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:471:CYS:HB3	8:A:852:HOH:O	2.17	0.44
1:A:73:GLY:N	1:A:74:PRO:CD	2.79	0.44
1:A:245:MET:HG3	1:A:324:CYS:HA	1.99	0.44
1:A:275:CYS:HA	1:A:358:LEU:HD22	2.00	0.44
1:A:22:GLU:OE2	1:A:26:LYS:HE3	2.17	0.44
2:D:96:VAL:HG21	2:D:115:SER:HB2	1.98	0.44
2:D:151:THR:HG23	2:D:162:LEU:HD11	1.99	0.44
1:C:378:GLY:HA3	1:C:401:TYR:CD1	2.52	0.44
1:C:465:MET:HG3	1:C:466:THR:N	2.32	0.44
2:D:143:LYS:N	2:D:144:PRO:HD3	2.31	0.44
1:A:65:ALA:O	1:A:70:VAL:HG13	2.17	0.44
2:B:93:GLN:HG2	8:B:780:HOH:O	2.18	0.44
1:C:441:MET:HB3	1:C:441:MET:HE2	1.77	0.44
2:B:82:PHE:C	2:B:85:THR:HG23	2.39	0.43
1:A:27:ASP:HA	1:A:30:LYS:HE3	2.00	0.43
2:B:279:GLU:OE1	2:B:279:GLU:C	2.57	0.43
1:C:106:VAL:HG22	1:C:107:ASN:N	2.34	0.43
1:C:382:ALA:HB1	1:C:386:ASP:HB2	1.99	0.43
2:D:132:LYS:HD3	2:D:174:PHE:CE2	2.53	0.43
2:D:323:ASP:HB2	2:D:381:LYS:NZ	2.33	0.43
1:C:35:ASN:ND2	1:C:391:MET:HA	2.33	0.43
1:C:57:MET:HB3	2:D:142:TYR:OH	2.18	0.43
2:D:7:LYS:NZ	2:D:8:ILE:O	2.51	0.43
1:A:162:ASP:OD2	1:A:165:SER:HB2	2.18	0.43
1:A:373:GLU:HG2	1:A:375:VAL:HG13	2.00	0.43
1:C:72:TRP:CZ2	1:C:202:VAL:HG22	2.52	0.43
1:C:347:GLU:OE1	1:C:371:GLY:O	2.37	0.43
1:A:454:ASP:OD2	2:B:2:SER:N	2.52	0.43
2:B:373:PRO:HD2	8:B:719:HOH:O	2.17	0.43
2:D:88:TYR:OH	2:D:116:ASP:HB3	2.18	0.43
2:D:254:LEU:O	2:D:255:SER:HB3	2.19	0.43
2:D:390:PRO:HB2	2:D:393:ILE:HD11	2.01	0.43
2:B:82:PHE:CB	2:B:85:THR:HG21	2.49	0.43
2:D:341:PRO:HD2	2:D:344:LEU:HD12	2.01	0.43
1:C:428:LYS:HE2	1:C:429:PHE:CE1	2.53	0.43
2:D:224:ILE:O	2:D:251:TYR:HA	2.19	0.42
2:B:108:ARG:O	2:B:109:GLU:HG2	2.19	0.42
2:B:369:LEU:HD22	2:B:369:LEU:C	2.40	0.42
1:C:5:SER:O	1:C:8:GLU:N	2.53	0.42
1:C:104:THR:HA	1:C:108:ALA:O	2.19	0.42
2:D:242:ARG:HD2	2:D:246:GLU:OE1	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:260:SER:OG	2:D:31:PHE:O	2.37	0.42
2:B:326:ASP:OD2	2:B:348:ARG:NH1	2.51	0.42
2:B:487:TYR:O	2:B:491:MET:HG3	2.20	0.42
1:C:59:ILE:HG12	1:C:427:GLU:OE2	2.20	0.42
1:C:106:VAL:HG21	2:D:40:ILE:HG23	2.01	0.42
1:A:387:TYR:O	1:A:391:MET:HG2	2.19	0.42
1:A:70:VAL:HG22	1:A:71:VAL:HG23	2.01	0.42
2:B:381:LYS:O	2:B:385:GLU:HG3	2.19	0.42
1:A:162:ASP:OD2	1:A:165:SER:HB3	2.19	0.42
1:A:443:SER:O	1:A:444:TRP:HB2	2.20	0.42
1:A:475:LEU:HD12	2:B:267:GLY:H	1.84	0.42
2:B:369:LEU:HD13	2:B:369:LEU:N	2.35	0.42
1:C:139:GLU:HG3	1:C:174:LEU:HD13	2.02	0.42
1:C:426:LYS:HB2	1:C:427:GLU:OE2	2.20	0.42
1:A:96:ARG:NH1	5:A:603:ICS:S5A	2.93	0.42
1:A:437:PRO:HG2	1:A:467:LEU:HD12	2.01	0.42
1:A:22:GLU:OE1	1:A:25:ARG:NH2	2.53	0.41
2:B:87:PRO:HG2	2:B:112:SER:O	2.19	0.41
2:D:300:LYS:HE2	8:D:792:HOH:O	2.19	0.41
1:A:345:ARG:HD2	8:A:883:HOH:O	2.19	0.41
2:D:86:MET:HG2	2:D:138:CYS:SG	2.60	0.41
2:D:478:HIS:O	2:D:481:ARG:HG3	2.20	0.41
1:C:253:TRP:HE1	1:C:265:THR:HG1	1.67	0.41
2:D:153:CYS:HB3	2:D:188:ALA:HB3	2.03	0.41
1:A:226:ILE:HA	1:A:253:TRP:HB2	2.03	0.41
2:D:194:VAL:HB	2:D:297:HIS:CG	2.55	0.41
1:A:365:GLY:CA	8:A:711:HOH:O	2.69	0.41
2:D:369:LEU:HD13	2:D:392:HIS:O	2.20	0.41
1:A:22:GLU:OE1	1:A:25:ARG:NH1	2.53	0.41
1:A:139:GLU:CD	1:A:176:LYS:HD2	2.41	0.41
1:A:187:ARG:HD3	8:B:975:HOH:O	2.20	0.41
1:A:361:ARG:HB3	1:A:379:TYR:OH	2.20	0.41
1:C:442:HIS:HB3	5:C:603:ICS:S4B	2.60	0.41
2:D:147:ILE:HG13	2:D:179:PHE:CE2	2.54	0.41
2:D:518:ASN:O	2:D:518:ASN:CG	2.58	0.41
2:B:143:LYS:HA	2:B:143:LYS:CE	2.41	0.41
2:D:23:MET:CE	2:D:24:LEU:HG	2.51	0.41
2:D:401:ARG:CZ	2:D:401:ARG:CB	2.99	0.41
1:A:50:LYS:NZ	8:A:731:HOH:O	2.51	0.41
1:A:83:HIS:HA	1:A:116:SER:OG	2.20	0.41
2:B:228:PRO:HA	2:B:293:LEU:HD12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:C:786:HOH:O	2:D:34:LYS:HE2	2.21	0.41
2:D:297:HIS:CD2	2:D:297:HIS:O	2.74	0.41
2:B:326:ASP:OD1	2:B:348:ARG:HD2	2.21	0.41
1:A:419:ASP:O	1:A:467:LEU:HD11	2.21	0.40
2:B:322:LEU:HD23	1:C:474:LYS:HG3	2.02	0.40
1:C:73:GLY:N	1:C:74:PRO:CD	2.84	0.40
2:B:509:THR:O	2:B:516:ASP:HA	2.22	0.40
1:C:387:TYR:O	1:C:391:MET:HG2	2.21	0.40
2:D:172:GLU:OE2	8:D:701:HOH:O	2.22	0.40
1:A:58:THR:OG1	1:A:403:ASP:OD1	2.31	0.40
2:D:326:ASP:O	2:D:330:MET:HG3	2.22	0.40
1:A:134:LEU:HD12	1:A:134:LEU:O	2.21	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	469/492 (95%)	446 (95%)	21 (4%)	2 (0%)	34	30
1	C	465/492 (94%)	447 (96%)	16 (3%)	2 (0%)	34	30
2	B	520/523 (99%)	500 (96%)	19 (4%)	1 (0%)	47	44
2	D	520/523 (99%)	503 (97%)	16 (3%)	1 (0%)	47	44
All	All	1974/2030 (97%)	1896 (96%)	72 (4%)	6 (0%)	41	37

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	6	ARG
2	B	255	SER
1	C	6	ARG

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Mol	Chain	Res	Type
2	D	255	SER
1	A	355	ILE
1	C	355	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	390/415 (94%)	372 (95%)	18 (5%)	27	23
1	C	384/415 (92%)	363 (94%)	21 (6%)	21	17
2	B	444/454 (98%)	432 (97%)	12 (3%)	44	46
2	D	441/454 (97%)	431 (98%)	10 (2%)	50	53
All	All	1659/1738 (96%)	1598 (96%)	61 (4%)	34	32

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	GLU
1	A	14	GLN
1	A	50	LYS
1	A	98	ASN
1	A	116	SER
1	A	134	LEU
1	A	150	VAL
1	A	182	ARG
1	A	199	ASN
1	A	213	ASP
1	A	277	ARG
1	A	288	GLU
1	A	401	TYR
1	A	409	PHE
1	A	445	ASP
1	A	451	HIS
1	A	475	LEU
1	A	480	GLU

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Mol	Chain	Res	Type
2	B	13	PRO
2	B	85	THR
2	B	154	MET
2	B	172	GLU
2	B	214	ASP
2	B	215	ASP
2	B	258	GLU
2	B	260	VAL
2	B	281	MET
2	B	369	LEU
2	B	401	ARG
2	B	431	ARG
1	C	6	ARG
1	C	12	LEU
1	C	14	GLN
1	C	17	LEU
1	C	70	VAL
1	C	98	ASN
1	C	134	LEU
1	C	150	VAL
1	C	168	LYS
1	C	182	ARG
1	C	223	VAL
1	C	264	LEU
1	C	318	GLU
1	C	347	GLU
1	C	373	GLU
1	C	384	ASN
1	C	401	TYR
1	C	409	PHE
1	C	445	ASP
1	C	475	LEU
1	C	476	GLN
2	D	2	SER
2	D	38	ASP
2	D	154	MET
2	D	177	ASP
2	D	202	GLU
2	D	206	ARG
2	D	212	SER
2	D	258	GLU
2	D	260	VAL

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Mol	Chain	Res	Type
2	D	369	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	ASN
2	B	58	GLN
1	C	384	ASN
1	C	476	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](#)

Of 10 ligands modelled in this entry, 4 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	UFF	A	601	2,1	0,21,21	-	-	-		
4	HCA	C	602	-	13,13,13	1.17	1 (7%)	14,18,18	1.45	2 (14%)
4	HCA	A	602	-	13,13,13	1.02	0	14,18,18	2.44	6 (42%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	ICS	A	603	1	18,30,30	3.73	10 (55%)	-		
5	ICS	C	603	1	18,30,30	3.07	11 (61%)	-		
3	UFF	C	601	2,1	0,21,21	-	-	-		

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	UFF	A	601	2,1	-	-	0/9/8/8
4	HCA	C	602	-	-	7/17/17/17	-
4	HCA	A	602	-	-	6/17/17/17	-
3	UFF	C	601	2,1	-	-	0/9/8/8

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	603	ICS	S4B-FE5	-8.19	2.12	2.32
5	A	603	ICS	S3B-FE7	-7.63	2.13	2.32
5	C	603	ICS	S4B-FE7	-6.63	2.16	2.32
5	A	603	ICS	S4B-FE7	-6.14	2.17	2.32
5	C	603	ICS	S3B-FE7	-6.05	2.17	2.32
5	A	603	ICS	S1B-FE5	-4.94	2.20	2.32
5	C	603	ICS	S1B-FE6	-3.81	2.23	2.32
5	C	603	ICS	S4B-FE5	-3.43	2.23	2.32
5	A	603	ICS	S3A-FE5	3.28	2.32	2.24
5	C	603	ICS	S4A-FE3	-3.25	2.24	2.32
5	A	603	ICS	S1B-FE6	-3.22	2.24	2.32
5	C	603	ICS	S1B-FE5	-3.12	2.24	2.32
5	A	603	ICS	S3B-FE6	-3.09	2.24	2.32
5	C	603	ICS	S3B-FE6	-2.94	2.25	2.32
5	C	603	ICS	S2A-FE2	-2.76	2.25	2.32
5	C	603	ICS	S2B-FE2	-2.75	2.18	2.24
5	A	603	ICS	S1A-FE2	-2.69	2.25	2.32
5	A	603	ICS	S2A-FE2	-2.46	2.26	2.32
5	C	603	ICS	S3A-FE5	2.40	2.30	2.24
5	A	603	ICS	S2B-FE6	-2.37	2.19	2.24
5	C	603	ICS	S1A-FE2	-2.12	2.27	2.32
4	C	602	HCA	O4-C6	-2.07	1.23	1.30

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	602	HCA	O6-C7-C3	5.34	122.32	113.05
4	A	602	HCA	O7-C3-C7	3.93	114.39	108.86
4	C	602	HCA	O7-C3-C7	3.30	113.50	108.86
4	A	602	HCA	O5-C7-C3	-3.29	117.60	122.25
4	C	602	HCA	O6-C7-C3	2.92	118.12	113.05
4	A	602	HCA	O3-C6-C5	-2.76	114.22	123.08
4	A	602	HCA	O4-C6-C5	2.61	122.42	114.03
4	A	602	HCA	O1-C1-C2	-2.17	116.61	122.94

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	602	HCA	C2-C3-C7-O6
4	A	602	HCA	C4-C3-C7-O5
4	A	602	HCA	C4-C3-C7-O6
4	C	602	HCA	C2-C3-C7-O6
4	C	602	HCA	C4-C3-C7-O6
4	C	602	HCA	O7-C3-C7-O6
4	A	602	HCA	C2-C3-C7-O5
4	C	602	HCA	C2-C3-C7-O5
4	C	602	HCA	C4-C3-C7-O5
4	C	602	HCA	C1-C2-C3-C4
4	A	602	HCA	C4-C5-C6-O4
4	A	602	HCA	C4-C5-C6-O3
4	C	602	HCA	C1-C2-C3-C7

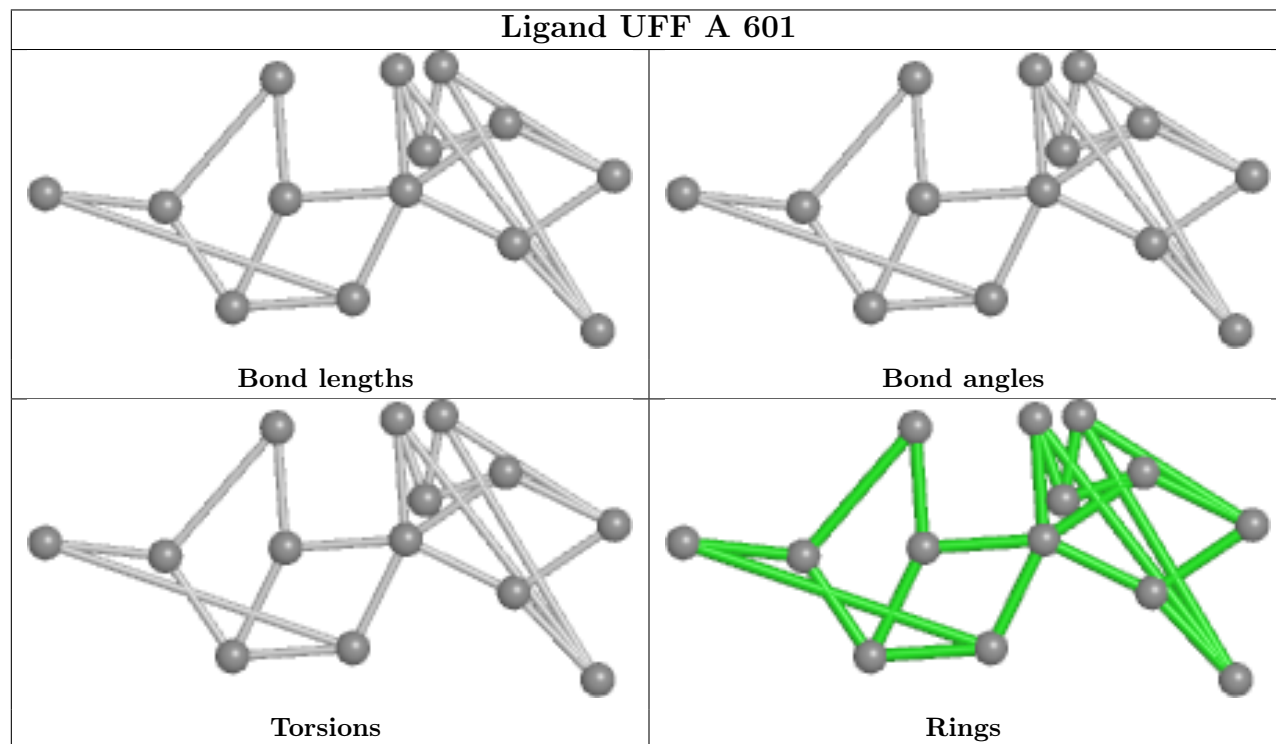
There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	UFF	4	0
5	A	603	ICS	1	0
5	C	603	ICS	2	0

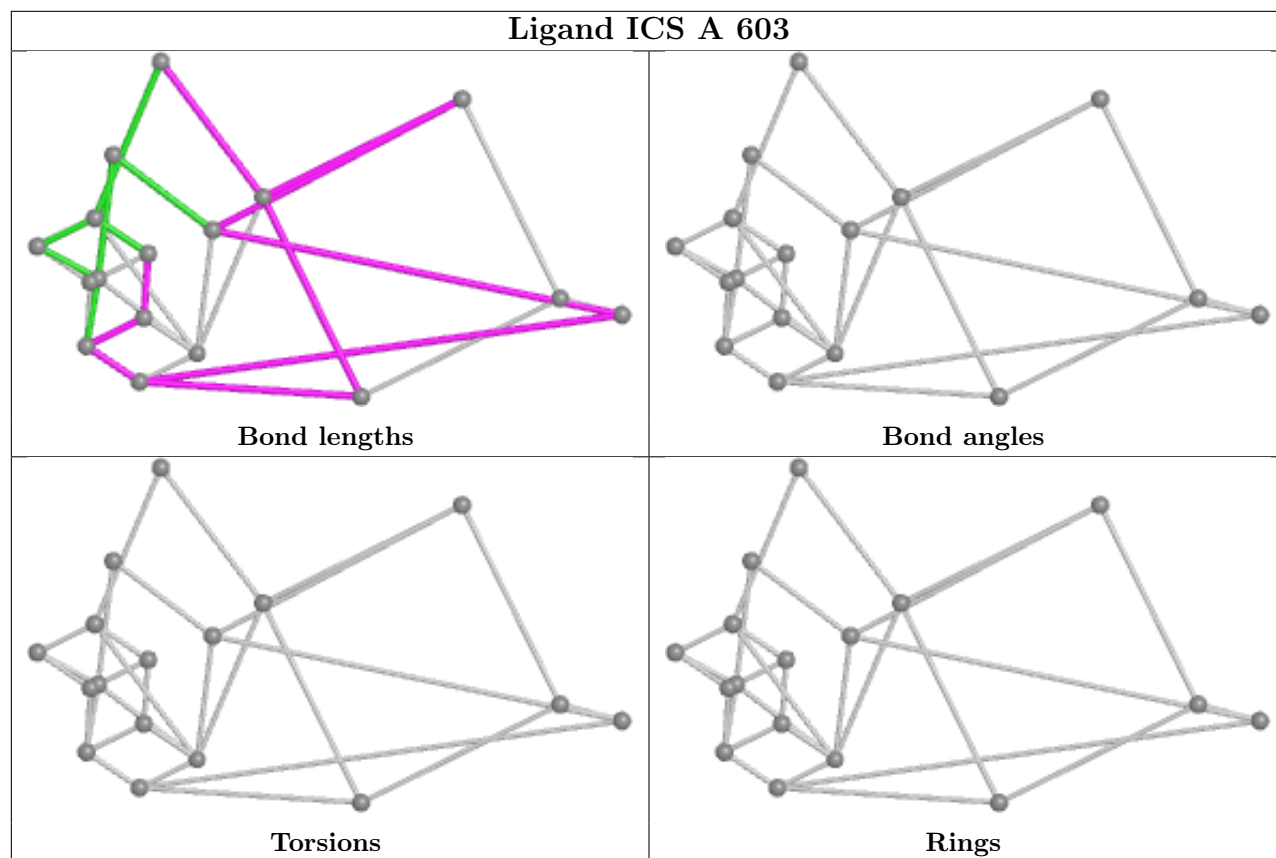
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be

highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

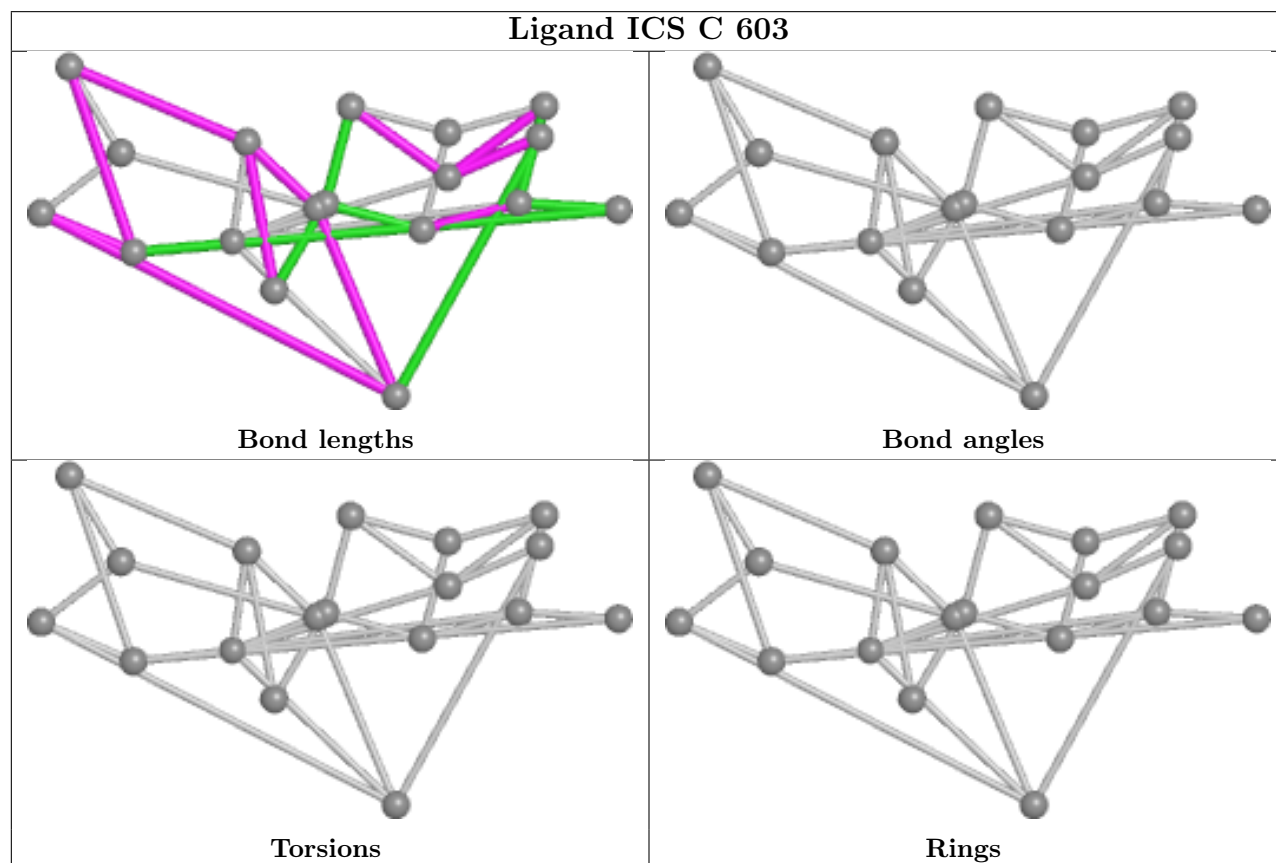


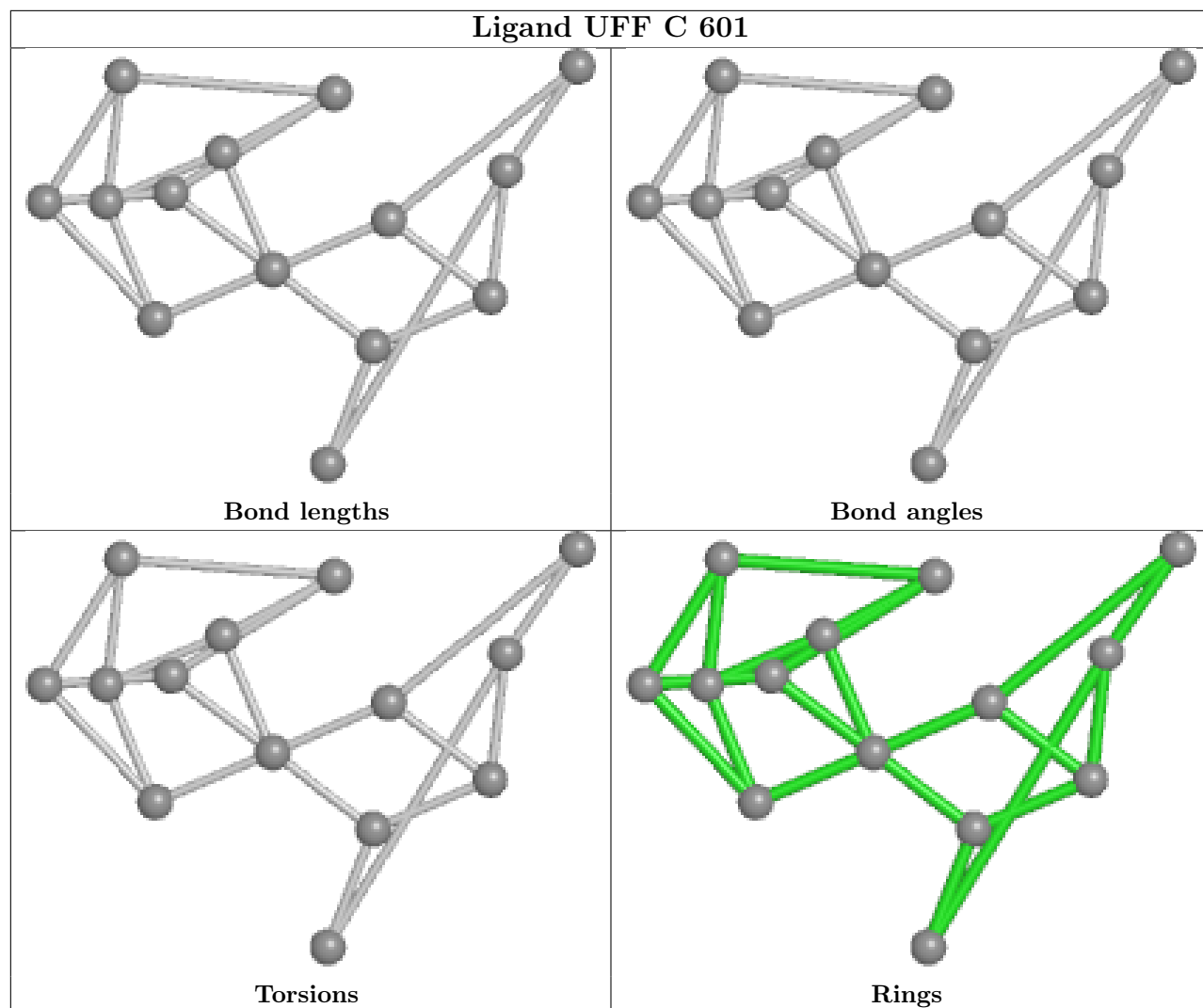


## Ligand ICS A 603



## Ligand ICS C 603





## 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 ⓘ

### 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<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	473/492 (96%)	0.12	10 (2%) 63 62	22, 37, 64, 87	0
1	C	469/492 (95%)	0.11	9 (1%) 66 65	22, 38, 60, 87	0
2	B	522/523 (99%)	-0.04	2 (0%) 92 92	18, 31, 45, 72	0
2	D	522/523 (99%)	-0.12	2 (0%) 92 92	21, 32, 46, 67	0
All	All	1986/2030 (97%)	0.01	23 (1%) 79 78	18, 34, 55, 87	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	40	THR	4.0
2	D	123	ALA	3.6
1	C	9	VAL	3.5
2	B	125	PHE	3.2
1	A	36	ASP	3.1
1	C	34	VAL	3.0
1	C	14	GLN	2.8
1	A	9	VAL	2.6
1	C	12	LEU	2.5
1	A	43	LYS	2.4
1	C	5	SER	2.4
1	C	124	VAL	2.4
1	A	400	LEU	2.4
1	C	253	TRP	2.4
1	A	22	GLU	2.3
1	A	29	ASN	2.3
1	A	46	ILE	2.2
2	D	125	PHE	2.2
1	C	134	LEU	2.1
1	C	35	ASN	2.1
2	B	122	ALA	2.1

*Continued on next page...*

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Mol	Chain	Res	Type	RSRZ
1	A	212	GLU	2.0
1	A	253	TRP	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

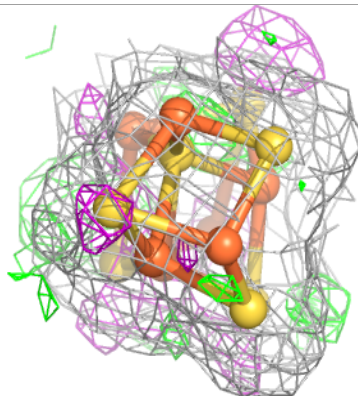
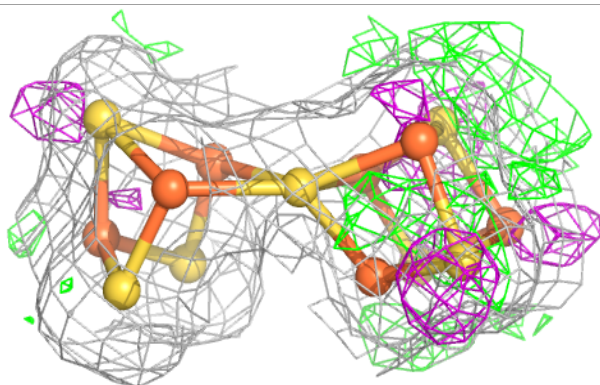
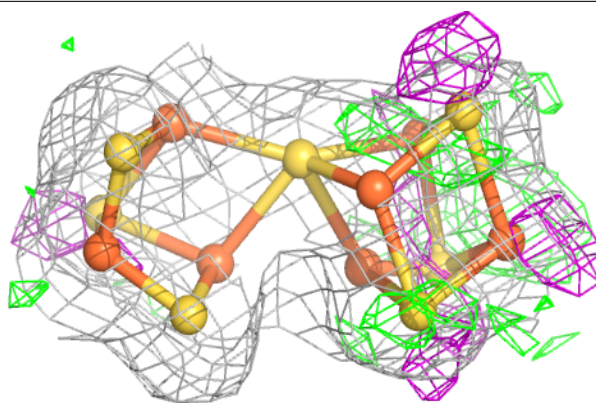
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
7	NA	B	602	1/1	0.81	0.08	25,25,25,25	1
7	NA	D	602	1/1	0.83	0.08	27,27,27,27	1
4	HCA	A	602	14/14	0.86	0.17	26,41,56,56	0
4	HCA	C	602	14/14	0.90	0.14	26,48,66,66	0
3	UFF	A	601	14/14	0.93	0.11	24,31,34,40	5
3	UFF	C	601	14/14	0.96	0.09	28,33,36,40	5
6	FE	B	601	1/1	0.98	0.09	32,32,32,32	1
6	FE	D	601	1/1	0.98	0.10	36,36,36,36	1
5	ICS	A	603	18/18	0.98	0.07	24,29,34,39	0
5	ICS	C	603	18/18	0.98	0.06	24,31,41,41	0

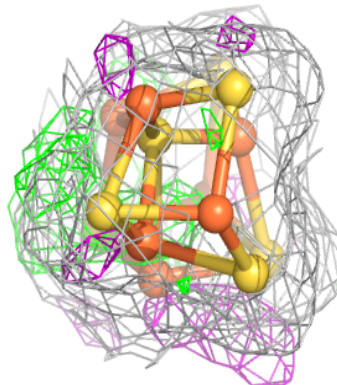
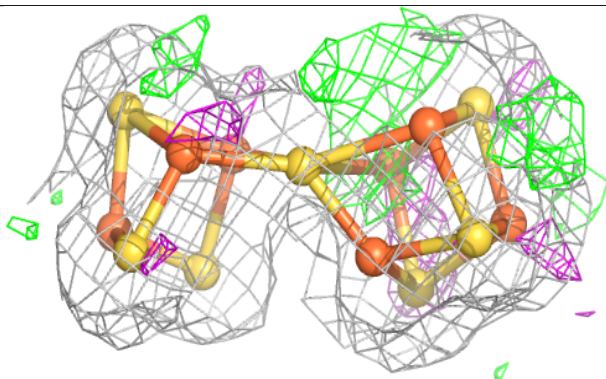
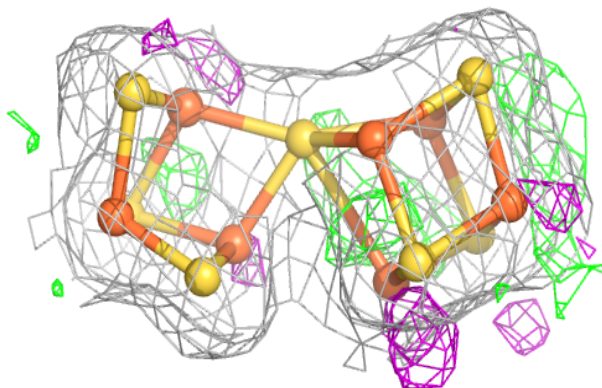
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around UFF A 601:**

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

**Electron density around UFF C 601:**

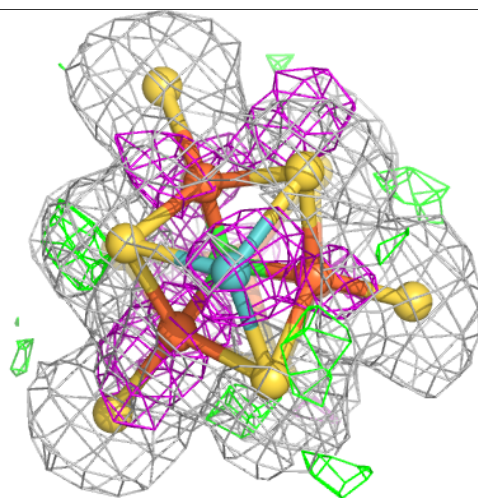
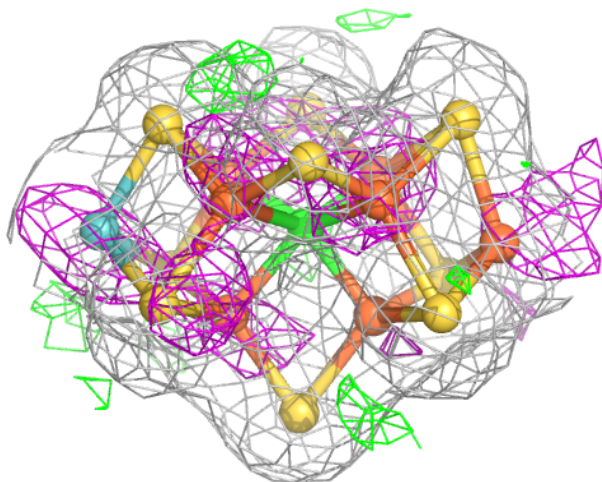
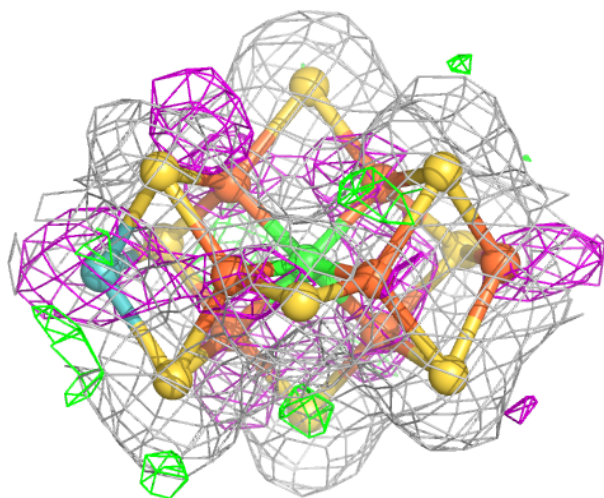
$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





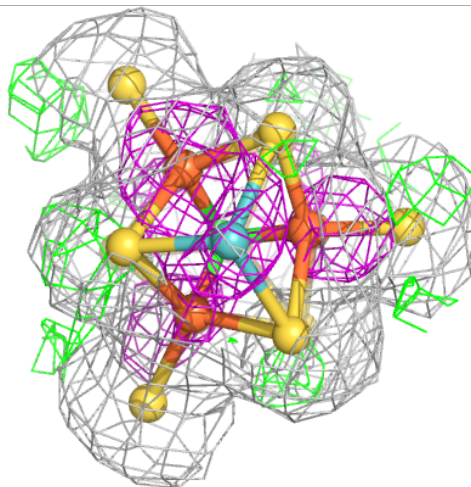
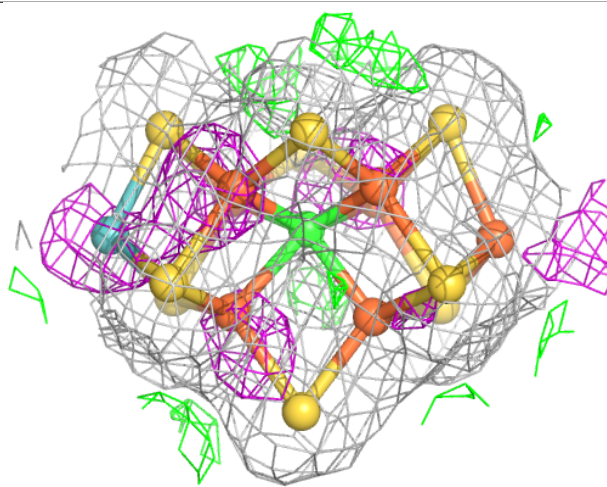
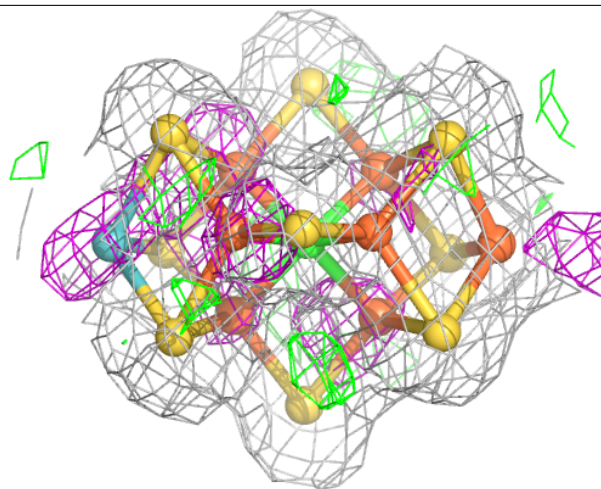
**Electron density around ICS A 603:**

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



**Electron density around ICS C 603:**

$2mF_o-DF_c$  (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.