



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 11:54 am BST

PDB ID : 4XPI
Title : Fe protein independent substrate reduction by nitrogenase variants altered in intramolecular electron transfer
Authors : Danyal, K.; Rasmusen, A.J.; Keable, S.M.; Shaw, S.; Zadvornyy, O.; Duval, S.; Dean, D.R.; Rauegi, S.; Peters, J.W.; Seefeldt, L.C.
Deposited on : 2015-01-17
Resolution : 1.97 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 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 : 2.11
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) : 2.11

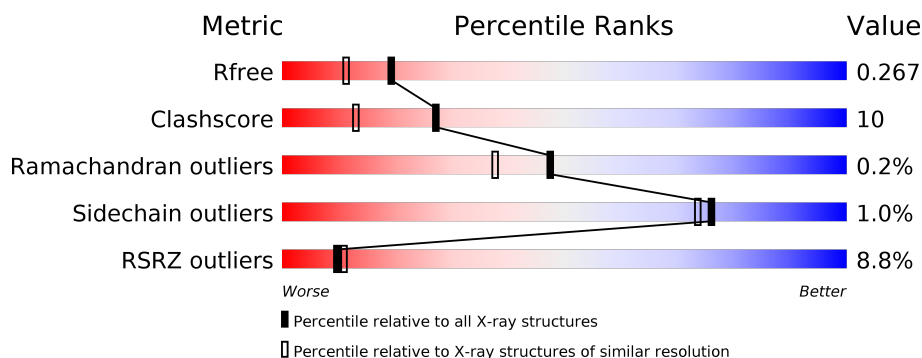
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.97 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	11647 (2.00-1.96)
Clashscore	141614	1014 (1.98-1.98)
Ramachandran outliers	138981	1006 (1.98-1.98)
Sidechain outliers	138945	1006 (1.98-1.98)
RSRZ outliers	127900	11410 (2.00-1.96)

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 on 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	490	<div> <div>11%</div> <div>79%</div> <div>18%</div> <div>..</div> </div>
1	C	490	<div> <div>16%</div> <div>71%</div> <div>25%</div> <div>..</div> </div>
2	B	522	<div> <div>4%</div> <div>86%</div> <div>13%</div> <div>.</div> </div>
2	D	522	<div> <div>4%</div> <div>87%</div> <div>13%</div> <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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	GOL	B	607	-	-	X	-
5	GOL	B	608	-	X	X	-
5	GOL	C	505	-	-	X	-
5	GOL	C	508	-	-	X	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 17075 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 molybdenum-iron protein alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	479	Total	C	N	O	S	1	2	0
			3805	2419	648	712	26			
1	C	478	Total	C	N	O	S	0	1	0
			3797	2414	647	711	25			

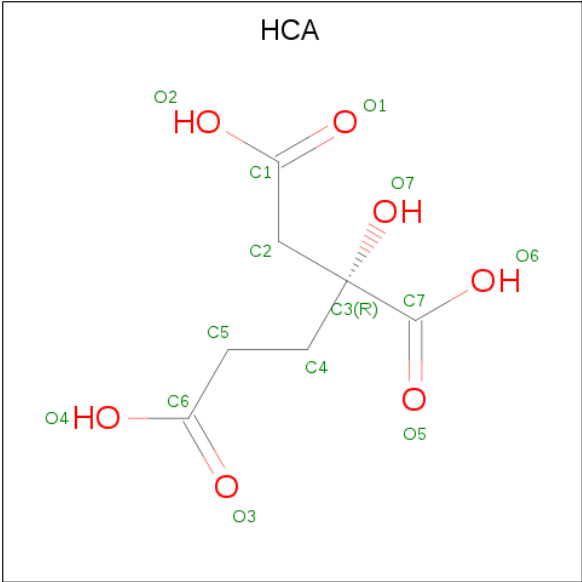
- Molecule 2 is a protein called Nitrogenase molybdenum-iron protein beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	522	Total	C	N	O	S	0	1	0
			4178	2666	708	776	28			
2	D	522	Total	C	N	O	S	0	0	0
			4172	2663	707	774	28			

There are 2 discrepancies between the modelled and reference sequences:

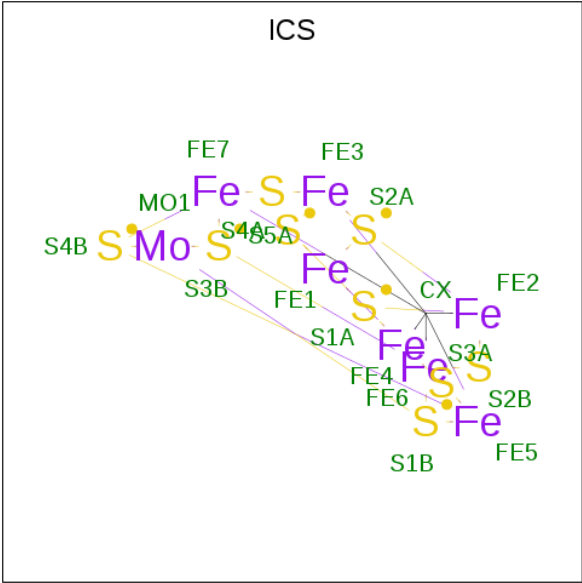
Chain	Residue	Modelled	Actual	Comment	Reference
B	98	HIS	TYR	engineered mutation	UNP P07329
D	98	HIS	TYR	engineered mutation	UNP P07329

- Molecule 3 is 3-HYDROXY-3-CARBOXY-ADIPIC ACID (three-letter code: HCA) (formula: C₇H₁₀O₇).



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 (three-letter code: ICS) (formula: CFe₇MoS₉).



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

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

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



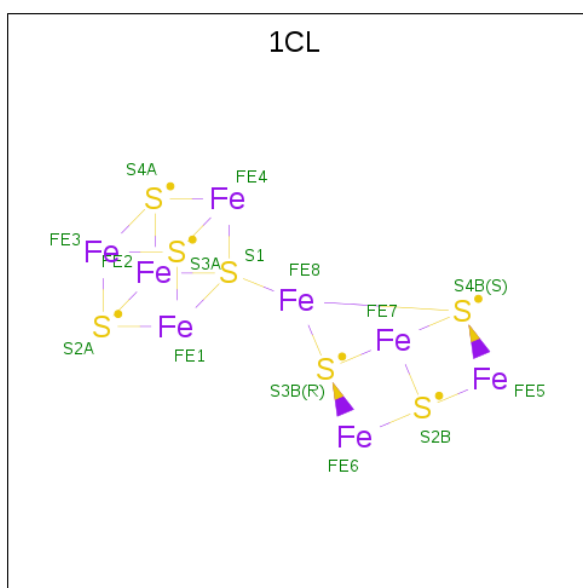
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	D	1	Total	C	O	0	0
			6	3	3		
5	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is FE(8)-S(7) CLUSTER, OXIDIZED (three-letter code: 1CL) (formula: Fe₈S₇).



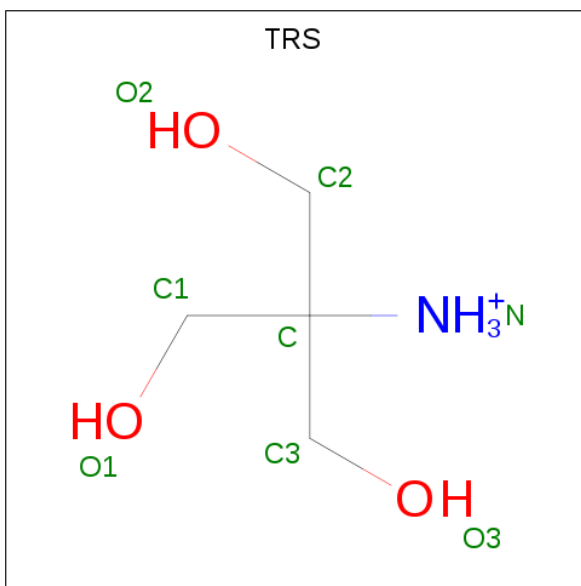
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	Fe	S	0	0
			15	8	7		
6	C	1	Total	Fe	S	0	0
			15	8	7		

- Molecule 7 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	2	Total	Ca	0	0
			2	2		

- Molecule 8 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code:

TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	B	1	Total	C	N	O	0	0
			8	4	1	3		
8	D	1	Total	C	N	O	0	0
			8	4	1	3		

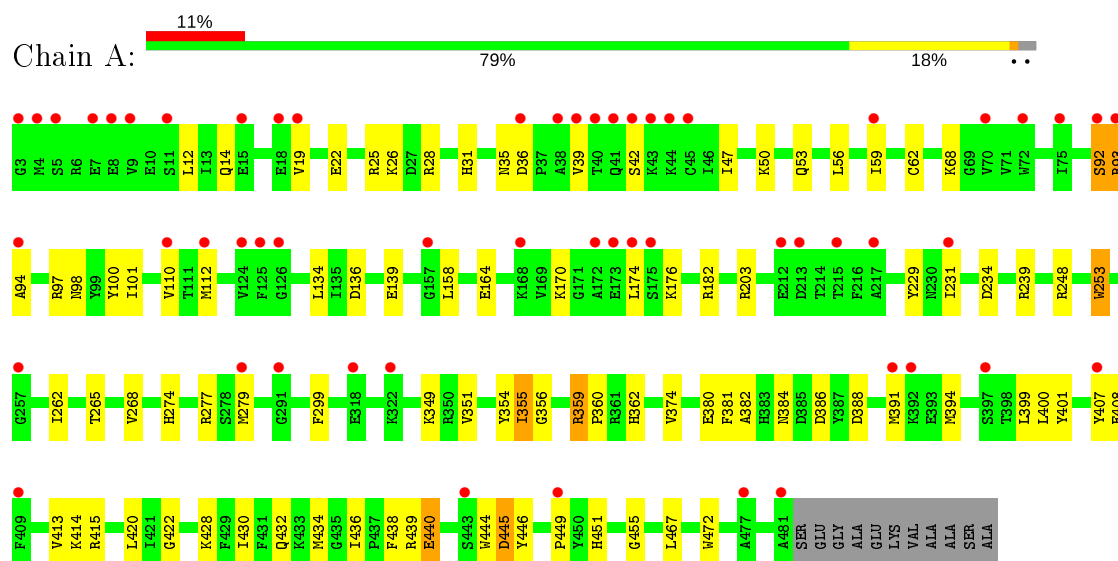
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	211	Total	O	0	0
			211	211		
9	B	257	Total	O	0	0
			257	257		
9	C	202	Total	O	0	0
			202	202		
9	D	251	Total	O	0	0
			251	251		

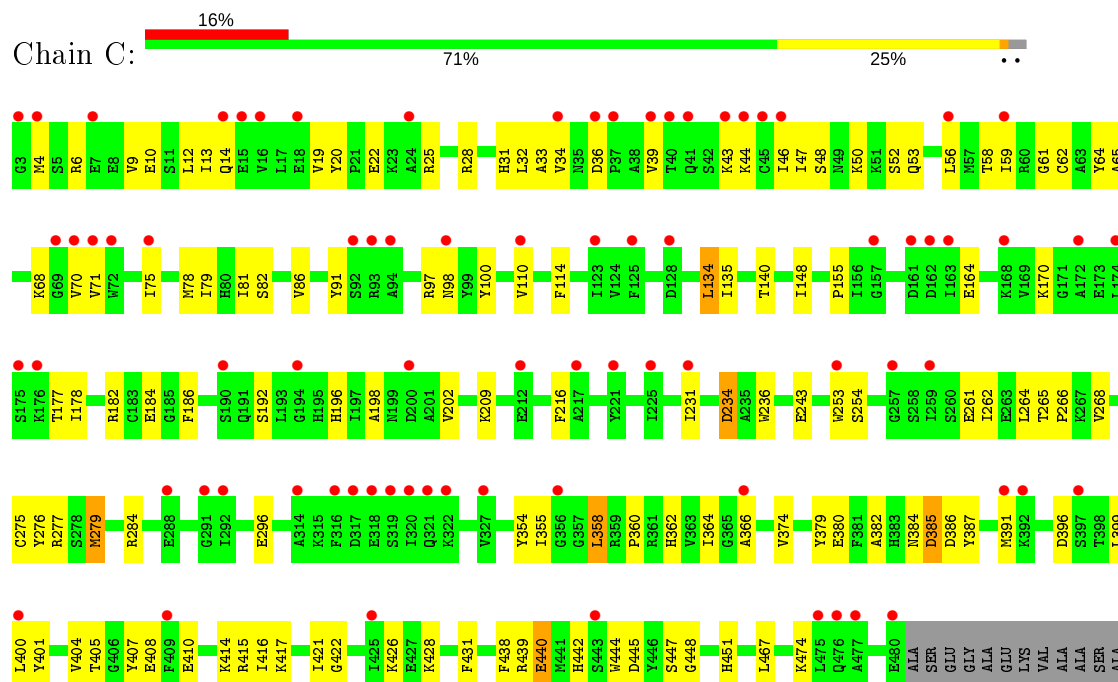
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

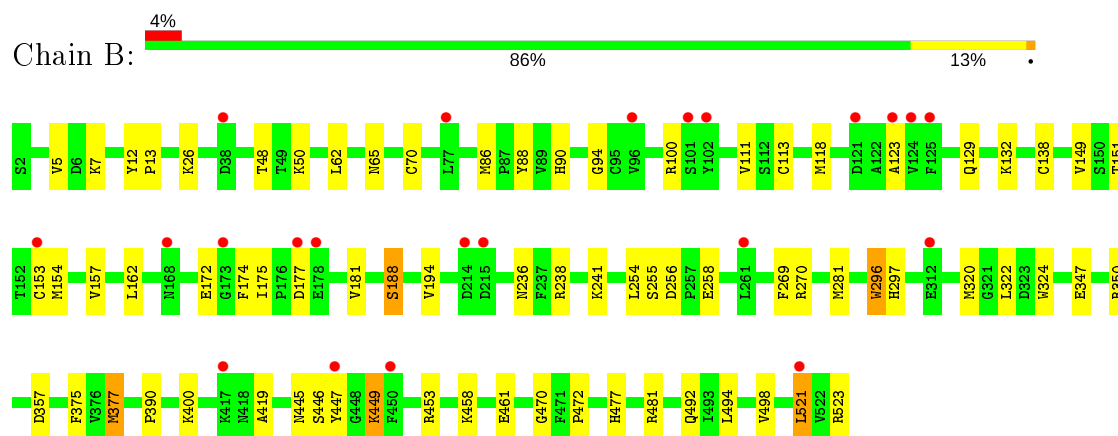
- Molecule 1: Nitrogenase molybdenum-iron protein alpha chain



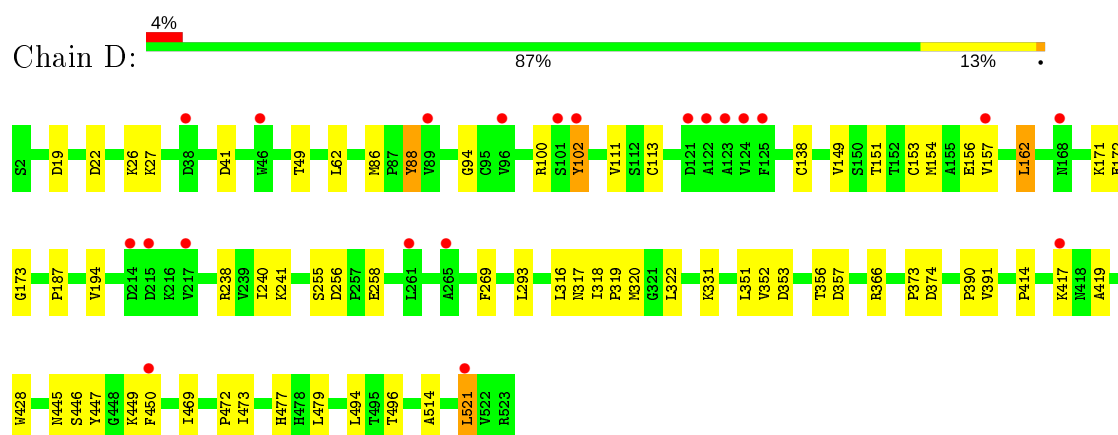
- Molecule 1: Nitrogenase molybdenum-iron protein alpha chain



- Molecule 2: Nitrogenase molybdenum-iron protein beta chain



- Molecule 2: Nitrogenase molybdenum-iron protein beta chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	80.80Å 130.83Å 108.11Å 90.00° 111.14° 90.00°	Depositor
Resolution (Å)	35.00 – 1.97 35.89 – 1.97	Depositor EDS
% Data completeness (in resolution range)	96.4 (35.00-1.97) 96.4 (35.89-1.97)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.96 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.213 , 0.264 0.218 , 0.267	Depositor DCC
R_{free} test set	7091 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	27.7	Xtriage
Anisotropy	0.531	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 49.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17075	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 24.75 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.6143e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: GOL, CA, HCA, ICS, TRS, 1CL

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	1.06	7/3899 (0.2%)	0.96	6/5257 (0.1%)
1	C	0.93	4/3888 (0.1%)	0.92	4/5242 (0.1%)
2	B	1.03	6/4284 (0.1%)	0.94	7/5791 (0.1%)
2	D	1.01	6/4278 (0.1%)	0.92	8/5783 (0.1%)
All	All	1.01	23/16349 (0.1%)	0.94	25/22073 (0.1%)

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	440	GLU	CD-OE2	-11.81	1.12	1.25
1	A	92	SER	CA-C	-11.61	1.22	1.52
1	C	440	GLU	CD-OE2	-8.55	1.16	1.25
2	B	357	ASP	C-O	7.68	1.38	1.23
1	C	91	TYR	CD1-CE1	-7.25	1.28	1.39
2	D	428	TRP	CD2-CE2	7.22	1.50	1.41
2	D	357	ASP	C-O	7.02	1.36	1.23
1	A	430	ILE	C-O	6.86	1.36	1.23
1	A	356	GLY	C-O	-6.78	1.12	1.23
2	B	188	SER	CB-OG	6.52	1.50	1.42
2	D	450	PHE	N-CA	6.23	1.58	1.46
2	D	477	HIS	CG-CD2	6.20	1.46	1.35
1	C	444	TRP	CD2-CE2	6.17	1.48	1.41
2	B	324	TRP	CD2-CE2	5.86	1.48	1.41
1	A	444	TRP	CD2-CE2	5.86	1.48	1.41
2	B	477	HIS	CG-CD2	5.68	1.45	1.35
2	D	102	TYR	CE2-CZ	-5.62	1.31	1.38
1	A	229	TYR	CG-CD2	5.48	1.46	1.39
2	D	496	THR	C-O	5.39	1.33	1.23
1	C	236	TRP	CD2-CE2	5.36	1.47	1.41
2	B	477	HIS	CE1-NE2	5.26	1.44	1.32
1	A	253	TRP	CG-CD1	5.12	1.44	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	296	TRP	CD2-CE2	5.00	1.47	1.41

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	320	MET	CG-SD-CE	7.86	112.78	100.20
2	D	521	LEU	CB-CG-CD2	7.45	123.66	111.00
1	A	93	ARG	NE-CZ-NH2	-6.94	116.83	120.30
2	B	521	LEU	CB-CG-CD2	6.89	122.72	111.00
2	B	377	MET	CG-SD-CE	-6.65	89.56	100.20
1	C	134	LEU	CB-CG-CD1	-6.21	100.45	111.00
1	C	234	ASP	CB-CG-OD1	6.11	123.80	118.30
1	C	358	LEU	CB-CG-CD2	-6.08	100.66	111.00
2	D	41	ASP	CB-CG-OD1	5.99	123.69	118.30
2	B	481	ARG	NE-CZ-NH2	-5.95	117.33	120.30
1	A	203	ARG	NE-CZ-NH1	5.74	123.17	120.30
1	A	359	ARG	NE-CZ-NH1	5.74	123.17	120.30
1	A	239	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	C	385	ASP	CB-CG-OD1	5.59	123.33	118.30
2	D	238	ARG	NE-CZ-NH2	-5.50	117.55	120.30
2	B	498	VAL	CG1-CB-CG2	-5.50	102.10	110.90
2	D	162	LEU	CB-CG-CD1	-5.31	101.97	111.00
2	D	469	ILE	CG1-CB-CG2	-5.29	99.75	111.40
2	D	19	ASP	CB-CG-OD1	5.24	123.01	118.30
1	A	445	ASP	CB-CG-OD1	5.23	123.01	118.30
2	B	238	ARG	NE-CZ-NH2	-5.09	117.75	120.30
2	B	350	ARG	NE-CZ-NH2	-5.05	117.78	120.30
2	B	523	ARG	NE-CZ-NH2	-5.05	117.78	120.30
1	A	248	ARG	NE-CZ-NH2	-5.05	117.78	120.30
2	D	374	ASP	CB-CG-OD1	5.03	122.82	118.30

There are no chirality outliers.

There are no planarity outliers.

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	3805	0	3747	83	0
1	C	3797	0	3737	135	0
2	B	4178	0	4089	63	0
2	D	4172	0	4085	45	0
3	A	14	0	6	1	0
3	C	14	0	6	2	0
4	A	18	0	0	0	0
4	C	18	0	0	0	0
5	A	6	0	8	0	0
5	B	36	0	48	17	0
5	C	36	0	48	19	0
5	D	12	0	16	0	0
6	A	15	0	0	0	0
6	C	15	0	0	0	0
7	B	2	0	0	0	0
8	B	8	0	12	2	0
8	D	8	0	12	1	0
9	A	211	0	0	22	0
9	B	257	0	0	25	0
9	C	202	0	0	52	0
9	D	251	0	0	16	0
All	All	17075	0	15814	324	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 10.

All (324) 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:C:71:VAL:HG13	9:C:796:HOH:O	1.24	1.35
1:C:428:LYS:HB2	9:C:766:HOH:O	1.39	1.17
1:C:385:ASP:HB2	9:C:691:HOH:O	1.38	1.17
1:C:358:LEU:HB2	9:C:615:HOH:O	1.61	0.99
1:C:46:ILE:HD12	9:C:730:HOH:O	1.60	0.98
2:B:111:VAL:HG12	9:B:735:HOH:O	1.67	0.95
5:C:504:GOL:O3	9:C:601:HOH:O	1.81	0.93
1:C:428:LYS:HD2	9:C:609:HOH:O	1.69	0.91
1:C:36:ASP:O	1:C:39:VAL:HG22	1.69	0.91
2:D:318:ILE:HA	9:D:701:HOH:O	1.69	0.91
1:A:394:MET:HB2	9:A:719:HOH:O	1.72	0.90
2:D:111:VAL:HG12	9:D:727:HOH:O	1.72	0.89
2:B:269:PHE:CD2	9:B:819:HOH:O	2.26	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:269:PHE:CD2	9:D:871:HOH:O	2.27	0.88
1:C:438:PHE:HB2	9:C:667:HOH:O	1.75	0.87
1:C:25:ARG:NH1	9:C:602:HOH:O	2.08	0.86
5:C:504:GOL:C3	9:C:601:HOH:O	2.23	0.86
1:C:140:THR:HG21	5:C:507:GOL:H12	1.55	0.86
2:D:194:VAL:HG21	9:D:925:HOH:O	1.76	0.85
1:A:440:GLU:HG3	9:A:785:HOH:O	1.75	0.85
1:A:428:LYS:HB2	9:A:785:HOH:O	1.76	0.84
1:C:6:ARG:NH2	1:C:396:ASP:OD1	2.11	0.83
1:A:36:ASP:O	1:A:39:VAL:HG22	1.80	0.82
1:C:28:ARG:HD3	1:C:408:GLU:OE2	1.79	0.82
1:C:186:PHE:HB3	9:D:773:HOH:O	1.78	0.81
1:A:445:ASP:OD2	9:A:601:HOH:O	2.00	0.80
1:C:13:ILE:HG21	9:C:741:HOH:O	1.81	0.79
1:C:416:ILE:HG13	9:C:737:HOH:O	1.82	0.79
2:B:90:HIS:HB2	9:B:801:HOH:O	1.82	0.78
2:D:100:ARG:HG3	9:D:932:HOH:O	1.84	0.78
1:A:440:GLU:CG	9:A:785:HOH:O	2.27	0.78
2:B:461:GLU:OE1	9:B:701:HOH:O	2.01	0.77
1:A:139:GLU:HG3	1:A:174:LEU:HD13	1.69	0.75
1:C:78:MET:HE2	9:C:628:HOH:O	1.87	0.75
1:A:432:GLN:OE1	9:A:602:HOH:O	2.05	0.74
2:B:88:TYR:HE2	9:B:801:HOH:O	1.70	0.74
2:D:293:LEU:O	9:D:701:HOH:O	2.03	0.74
1:C:61:GLY:HA3	9:C:656:HOH:O	1.86	0.74
1:C:234:ASP:HB3	1:C:451:HIS:ND1	2.03	0.74
1:C:58:THR:HB	5:C:508:GOL:O1	1.89	0.72
1:C:64:TYR:HB2	9:C:627:HOH:O	1.88	0.72
1:C:253:TRP:HB3	1:C:279:MET:CE	2.20	0.72
1:C:385:ASP:CB	9:C:691:HOH:O	2.14	0.72
1:A:253:TRP:HB3	1:A:279:MET:HE1	1.72	0.70
2:B:129:GLN:HA	9:B:718:HOH:O	1.90	0.70
1:C:410:GLU:OE2	1:C:414:LYS:NZ	2.23	0.69
1:A:42:SER:N	1:A:388:ASP:OD1	2.19	0.69
1:A:53:GLN:HB2	1:A:56:LEU:HD12	1.73	0.69
5:C:504:GOL:H31	9:C:601:HOH:O	1.91	0.69
1:C:275:CYS:HB2	9:C:765:HOH:O	1.93	0.68
2:B:347:GLU:OE2	9:B:702:HOH:O	2.12	0.67
1:C:12:LEU:HD13	1:C:415:ARG:HG2	1.76	0.66
1:C:4:MET:SD	9:C:737:HOH:O	2.54	0.66
1:A:351:VAL:CG2	1:A:374:VAL:HG22	2.26	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:373:PRO:HD2	9:D:925:HOH:O	1.96	0.66
1:A:445:ASP:CG	9:A:601:HOH:O	2.34	0.66
1:A:428:LYS:NZ	9:A:605:HOH:O	2.28	0.66
1:C:467:LEU:O	1:C:467:LEU:HD23	1.96	0.66
1:C:28:ARG:CD	1:C:408:GLU:OE2	2.42	0.65
1:C:33:ALA:HB3	9:C:730:HOH:O	1.97	0.65
1:C:53:GLN:HB2	1:C:56:LEU:HD12	1.79	0.65
1:C:209:LYS:HE2	9:C:659:HOH:O	1.97	0.64
2:B:236:ASN:ND2	5:B:607:GOL:H12	2.13	0.64
1:A:93:ARG:NH2	5:B:608:GOL:O2	2.31	0.64
1:C:10:GLU:HG3	1:C:34:VAL:HG21	1.79	0.64
2:B:129:GLN:OE1	9:B:703:HOH:O	2.15	0.63
1:A:101:ILE:HD12	5:B:603:GOL:H31	1.80	0.63
1:C:416:ILE:CG1	9:C:737:HOH:O	2.42	0.63
2:D:353:ASP:OD2	9:D:702:HOH:O	2.16	0.63
1:A:62:CYS:HA	9:A:641:HOH:O	1.98	0.62
1:C:32:LEU:HD21	9:C:753:HOH:O	1.98	0.62
1:A:391:MET:HA	9:A:719:HOH:O	2.00	0.62
1:A:35:ASN:HB2	1:A:400:LEU:HD11	1.82	0.62
1:C:192:SER:OG	9:C:603:HOH:O	2.16	0.62
1:C:253:TRP:HB3	1:C:279:MET:HE1	1.82	0.61
1:A:134:LEU:HG	2:B:62:LEU:HB2	1.82	0.61
1:C:71:VAL:CG1	9:C:796:HOH:O	2.04	0.61
1:C:53:GLN:O	5:C:508:GOL:C1	2.49	0.61
1:C:440:GLU:HG3	9:C:766:HOH:O	2.00	0.60
1:C:53:GLN:N	5:C:508:GOL:H11	2.16	0.60
1:C:14:GLN:HB3	9:C:764:HOH:O	2.01	0.60
1:C:253:TRP:HB3	1:C:279:MET:HE3	1.82	0.60
1:C:400:LEU:CB	9:C:730:HOH:O	2.49	0.60
2:B:7:LYS:HD3	9:B:810:HOH:O	2.02	0.59
2:B:446:SER:O	2:B:449:LYS:HG2	2.02	0.59
2:B:65:ASN:HD21	5:B:608:GOL:H31	1.67	0.58
9:C:604:HOH:O	2:D:26:LYS:HB2	2.04	0.58
2:B:172:GLU:OE1	9:B:704:HOH:O	2.17	0.58
1:C:47:ILE:CD1	1:C:50:LYS:HG3	2.34	0.58
1:C:438:PHE:CB	9:C:667:HOH:O	2.41	0.58
2:D:366:ARG:NH2	2:D:391:VAL:HG21	2.18	0.57
1:C:400:LEU:HB3	9:C:730:HOH:O	2.03	0.57
1:C:254:SER:HB3	9:C:748:HOH:O	2.04	0.57
1:C:81:ILE:C	9:C:630:HOH:O	2.43	0.57
1:C:47:ILE:HD12	1:C:50:LYS:HG3	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:MET:HG2	5:B:608:GOL:O2	2.05	0.57
1:A:39:VAL:HG23	1:A:39:VAL:O	2.05	0.56
1:A:445:ASP:O	1:A:446:TYR:HB2	2.05	0.56
2:B:90:HIS:CE1	2:B:118:MET:SD	2.99	0.56
1:C:405:THR:HG22	5:C:508:GOL:O3	2.05	0.56
2:B:100:ARG:HG3	9:B:947:HOH:O	2.04	0.56
1:C:216:PHE:CD1	1:C:264:LEU:HD22	2.41	0.56
1:A:47:ILE:HD12	1:A:50:LYS:HG2	1.89	0.55
1:C:48:SER:OG	1:C:384:ASN:ND2	2.38	0.55
2:D:390:PRO:O	2:D:419:ALA:HB2	2.06	0.55
1:C:39:VAL:O	1:C:39:VAL:HG23	2.07	0.55
1:A:164:GLU:OE1	1:A:182:ARG:NH2	2.41	0.54
1:C:467:LEU:C	1:C:467:LEU:HD23	2.28	0.54
2:D:113:CYS:HB2	9:D:727:HOH:O	2.06	0.54
1:A:22:GLU:HG3	1:A:25:ARG:NH2	2.23	0.54
1:C:265:THR:O	1:C:268:VAL:HG22	2.06	0.54
1:C:68:LYS:HD3	1:C:68:LYS:C	2.28	0.54
1:A:354:TYR:OH	1:A:380:GLU:HA	2.08	0.54
2:B:151:THR:HG23	2:B:162:LEU:HD11	1.89	0.54
1:C:164:GLU:OE1	1:C:182:ARG:NH2	2.40	0.54
1:C:44:LYS:NZ	9:C:614:HOH:O	2.39	0.53
1:C:65:ALA:O	1:C:70:VAL:HG23	2.08	0.53
1:C:53:GLN:O	5:C:508:GOL:H12	2.08	0.53
1:C:59:ILE:HD12	1:C:354:TYR:CE2	2.43	0.53
1:A:440:GLU:HG3	9:A:601:HOH:O	2.08	0.53
1:C:110:VAL:O	5:C:505:GOL:H11	2.08	0.53
1:A:164:GLU:OE2	1:A:182:ARG:NE	2.41	0.53
1:A:28:ARG:HD3	1:A:408:GLU:OE2	2.08	0.53
2:B:7:LYS:HE2	9:B:946:HOH:O	2.09	0.53
1:A:139:GLU:OE2	1:A:176:LYS:HE2	2.10	0.52
1:A:349:LYS:HE2	9:A:630:HOH:O	2.09	0.52
1:A:384:ASN:ND2	9:A:607:HOH:O	2.40	0.52
1:A:428:LYS:HD2	9:A:601:HOH:O	2.10	0.52
1:C:198:ALA:O	1:C:202:VAL:HG23	2.09	0.52
1:C:354:TYR:OH	1:C:380:GLU:HA	2.08	0.52
2:D:153:CYS:O	2:D:157:VAL:HG23	2.10	0.52
1:C:177:THR:HG21	9:C:659:HOH:O	2.10	0.52
1:C:400:LEU:HB2	9:C:730:HOH:O	2.10	0.52
1:C:428:LYS:NZ	9:C:607:HOH:O	2.28	0.52
2:B:390:PRO:O	2:B:419:ALA:HB2	2.10	0.51
2:B:50:LYS:HB2	8:B:604:TRS:N	2.24	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:404:VAL:HG23	9:C:753:HOH:O	2.10	0.51
1:A:12:LEU:HD13	1:A:415:ARG:HG2	1.92	0.51
2:B:453:ARG:NH1	5:B:608:GOL:O1	2.43	0.51
2:D:445:ASN:HB2	2:D:472:PRO:O	2.10	0.51
1:A:14:GLN:NE2	9:A:623:HOH:O	2.43	0.51
1:C:253:TRP:CB	1:C:279:MET:HE3	2.39	0.51
1:C:32:LEU:CD2	9:C:753:HOH:O	2.57	0.51
2:B:322:LEU:HD21	1:C:474:LYS:HB3	1.92	0.51
1:C:355:ILE:HB	1:C:360:PRO:HD3	1.92	0.51
1:C:140:THR:CG2	5:C:507:GOL:H12	2.36	0.51
1:C:354:TYR:CZ	1:C:404:VAL:HG12	2.46	0.51
1:A:445:ASP:O	1:A:446:TYR:CB	2.57	0.50
1:A:432:GLN:HG2	1:A:472:TRP:HH2	1.76	0.50
1:A:274:HIS:HE1	1:A:299:PHE:H	1.58	0.50
1:A:265:THR:O	1:A:268:VAL:HG22	2.11	0.50
1:C:135:ILE:HD12	1:C:170:LYS:HG3	1.92	0.50
1:C:52:SER:HB3	5:C:508:GOL:H11	1.92	0.50
1:A:28:ARG:CD	1:A:408:GLU:OE2	2.60	0.50
2:B:132:LYS:HD2	9:B:718:HOH:O	2.10	0.50
3:C:501:HCA:O7	3:C:501:HCA:O1	2.28	0.50
2:B:445:ASN:HA	9:B:725:HOH:O	2.11	0.50
9:A:637:HOH:O	2:B:26:LYS:HB2	2.12	0.49
2:D:154:MET:SD	9:D:773:HOH:O	2.59	0.49
2:B:296:TRP:CD1	2:B:377:MET:CE	2.95	0.49
1:C:75:ILE:HD13	1:C:262:ILE:HD11	1.95	0.49
5:B:603:GOL:H2	9:B:875:HOH:O	2.12	0.49
1:C:110:VAL:O	5:C:505:GOL:C1	2.60	0.49
3:A:501:HCA:O7	3:A:501:HCA:O2	2.29	0.49
1:A:253:TRP:HB3	1:A:279:MET:CE	2.41	0.49
1:A:47:ILE:HD11	9:A:611:HOH:O	2.11	0.49
2:B:123:ALA:HB2	2:B:154:MET:CE	2.43	0.49
2:D:171:LYS:HE2	9:D:815:HOH:O	2.13	0.49
2:B:153:CYS:O	2:B:157:VAL:HG23	2.13	0.48
1:C:19:VAL:HG11	1:C:407:TYR:CZ	2.48	0.48
1:C:75:ILE:HD11	9:C:748:HOH:O	2.12	0.48
1:C:243:GLU:OE1	9:C:604:HOH:O	2.20	0.48
1:A:413:VAL:CG1	1:A:436:ILE:CD1	2.91	0.48
2:D:319:PRO:HD3	9:D:701:HOH:O	2.12	0.48
1:A:62:CYS:HB3	2:B:94:GLY:HA3	1.93	0.48
1:C:53:GLN:O	5:C:508:GOL:H11	2.12	0.48
1:A:274:HIS:CE1	1:A:299:PHE:H	2.31	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:LYS:C	1:A:68:LYS:HD3	2.33	0.48
1:C:100:TYR:CE2	1:C:110:VAL:HB	2.48	0.48
2:B:132:LYS:CD	9:B:718:HOH:O	2.62	0.47
2:B:400:LYS:H	5:B:602:GOL:H32	1.79	0.47
1:A:100:TYR:CE2	1:A:110:VAL:HB	2.49	0.47
1:C:81:ILE:HD12	1:C:134:LEU:HD21	1.96	0.47
1:A:93:ARG:NH2	5:B:608:GOL:C2	2.77	0.47
2:D:473:ILE:HG21	2:D:479:LEU:HD12	1.97	0.47
1:C:442:HIS:HB3	3:C:501:HCA:O5	2.13	0.47
2:B:111:VAL:CG1	9:B:735:HOH:O	2.42	0.47
1:C:134:LEU:HG	2:D:62:LEU:HB2	1.97	0.47
1:C:405:THR:HG22	5:C:508:GOL:C3	2.45	0.47
1:A:428:LYS:HB2	1:A:438:PHE:CE2	2.50	0.47
2:B:296:TRP:CD1	2:B:377:MET:HE1	2.50	0.47
1:A:12:LEU:HD12	9:A:649:HOH:O	2.15	0.47
1:A:445:ASP:OD1	9:A:601:HOH:O	2.19	0.47
2:D:446:SER:O	2:D:449:LYS:HG2	2.14	0.47
1:A:422:GLY:HA2	1:A:439:ARG:O	2.15	0.47
1:C:31:HIS:NE2	1:C:47:ILE:HD11	2.30	0.47
1:C:79:ILE:HG23	1:C:114:PHE:CD2	2.50	0.47
2:D:173:GLY:HA2	9:D:791:HOH:O	2.14	0.46
1:A:351:VAL:HG21	1:A:374:VAL:HG22	1.97	0.46
1:C:261:GLU:CD	2:D:27:LYS:HZ3	2.19	0.46
1:A:47:ILE:HD12	1:A:50:LYS:CG	2.45	0.46
1:C:9:VAL:CG1	1:C:34:VAL:HG22	2.45	0.46
1:C:59:ILE:O	1:C:426:LYS:NZ	2.47	0.46
2:B:494:LEU:C	2:B:494:LEU:HD23	2.36	0.46
1:C:20:TYR:CE2	1:C:28:ARG:HG3	2.51	0.46
1:C:253:TRP:CZ2	1:C:262:ILE:HG23	2.50	0.46
1:C:431:PHE:HB2	9:C:667:HOH:O	2.14	0.46
1:C:62:CYS:HB3	2:D:94:GLY:HA3	1.97	0.46
1:C:155:PRO:HB3	2:D:157:VAL:HG21	1.97	0.46
1:C:192:SER:CB	9:C:603:HOH:O	2.63	0.46
2:B:88:TYR:O	2:B:149:VAL:HA	2.16	0.45
1:C:440:GLU:CG	9:C:766:HOH:O	2.61	0.45
1:C:428:LYS:HG3	1:C:438:PHE:CD2	2.51	0.45
1:A:234:ASP:HB3	1:A:451:HIS:ND1	2.30	0.45
1:C:277:ARG:HD2	1:C:386:ASP:OD2	2.16	0.45
1:C:417:LYS:NZ	9:C:605:HOH:O	2.22	0.45
1:A:414:LYS:HG3	1:A:434:MET:HE1	1.99	0.45
1:A:399:LEU:HD12	1:A:400:LEU:H	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:270:ARG:NH2	9:B:720:HOH:O	2.50	0.45
2:D:151:THR:CG2	2:D:162:LEU:HD21	2.46	0.45
2:D:22:ASP:O	2:D:26:LYS:HG2	2.17	0.45
1:A:59:ILE:HD12	1:A:354:TYR:CE2	2.52	0.44
2:B:70:CYS:HB2	2:B:188:SER:HB2	1.98	0.44
1:C:296:GLU:HB3	9:C:727:HOH:O	2.16	0.44
1:A:440:GLU:HG2	9:A:785:HOH:O	2.08	0.44
1:C:265:THR:HB	1:C:266:PRO:HD3	1.99	0.44
2:B:445:ASN:HB2	2:B:472:PRO:O	2.17	0.44
2:B:5:VAL:HG12	9:B:887:HOH:O	2.16	0.44
1:C:276:TYR:OH	1:C:284:ARG:NH2	2.50	0.44
2:D:86:MET:HG2	2:D:138:CYS:SG	2.57	0.44
1:A:93:ARG:HH21	5:B:608:GOL:HO2	1.63	0.44
1:C:421:ILE:O	1:C:438:PHE:HA	2.17	0.44
1:C:82:SER:N	9:C:630:HOH:O	2.50	0.44
2:D:322:LEU:HD13	2:D:351:LEU:HD23	1.99	0.44
2:B:7:LYS:CD	9:B:946:HOH:O	2.66	0.44
2:B:132:LYS:HD3	2:B:174:PHE:CE2	2.53	0.44
5:B:605:GOL:H12	9:B:829:HOH:O	2.17	0.44
1:C:382:ALA:HB3	1:C:387:TYR:CE2	2.53	0.44
1:A:381:PHE:HA	9:A:747:HOH:O	2.18	0.44
2:B:492:GLN:HB3	5:B:605:GOL:H31	2.00	0.44
1:C:6:ARG:HG2	1:C:10:GLU:OE2	2.18	0.44
2:B:194:VAL:HB	2:B:297:HIS:CG	2.53	0.43
2:D:390:PRO:O	2:D:419:ALA:CB	2.65	0.43
5:B:606:GOL:H12	5:B:607:GOL:O2	2.18	0.43
2:B:86:MET:HG2	2:B:138:CYS:SG	2.58	0.43
1:C:234:ASP:HB3	1:C:451:HIS:CG	2.54	0.43
1:C:79:ILE:O	1:C:148:ILE:HA	2.18	0.43
2:B:241:LYS:NZ	2:B:256:ASP:OD2	2.47	0.43
1:C:366:ALA:HB3	9:C:786:HOH:O	2.17	0.43
1:C:62:CYS:HA	9:C:644:HOH:O	2.18	0.43
1:C:9:VAL:HG12	1:C:34:VAL:HG22	2.01	0.43
2:D:49:THR:HB	8:D:602:TRS:C3	2.49	0.43
2:B:320:MET:SD	5:B:607:GOL:C3	3.07	0.43
2:B:90:HIS:CB	9:B:801:HOH:O	2.53	0.43
1:C:9:VAL:O	1:C:13:ILE:HD13	2.19	0.43
2:D:317:ASN:ND2	9:D:725:HOH:O	2.52	0.43
1:C:196:HIS:ND1	9:C:606:HOH:O	2.22	0.43
1:C:405:THR:HG22	5:C:508:GOL:H32	2.01	0.43
1:A:112:MET:HA	5:B:608:GOL:HO2	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:382:ALA:HB1	1:A:386:ASP:HB2	2.01	0.42
2:B:447:TYR:CD1	2:D:521:LEU:HD22	2.54	0.42
1:A:359:ARG:N	1:A:360:PRO:CD	2.82	0.42
2:B:12:TYR:HA	2:B:13:PRO:HA	1.83	0.42
1:C:184:GLU:HA	9:C:627:HOH:O	2.19	0.42
1:C:253:TRP:HA	1:C:254:SER:HA	1.75	0.42
2:D:352:VAL:O	2:D:356:THR:HG23	2.19	0.42
2:D:494:LEU:HD23	2:D:494:LEU:C	2.40	0.42
2:B:377:MET:HE2	2:B:377:MET:HB3	1.79	0.42
1:C:97:ARG:O	1:C:231:ILE:HA	2.19	0.42
2:D:156:GLU:HG3	2:D:187:PRO:HB3	2.01	0.42
1:A:158:LEU:HD11	2:B:154:MET:HG3	2.01	0.42
2:B:123:ALA:HB2	2:B:154:MET:HE3	2.02	0.42
1:C:364:ILE:HG12	1:C:374:VAL:HG21	2.01	0.42
2:B:470:GLY:N	9:B:725:HOH:O	2.52	0.42
2:D:240:ILE:HG13	2:D:293:LEU:HD11	2.01	0.42
1:A:399:LEU:HD12	1:A:400:LEU:N	2.35	0.42
2:B:50:LYS:HB2	8:B:604:TRS:HN1	1.84	0.42
1:C:399:LEU:HD12	1:C:400:LEU:H	1.85	0.42
1:C:86:VAL:O	1:C:86:VAL:HG22	2.19	0.42
2:D:154:MET:CG	9:D:773:HOH:O	2.68	0.42
1:A:420:LEU:HB2	1:A:467:LEU:HD12	2.01	0.41
1:A:413:VAL:HG11	1:A:436:ILE:HD12	2.02	0.41
1:A:449:PRO:O	1:A:455:GLY:HA2	2.19	0.41
2:B:470:GLY:CA	9:B:725:HOH:O	2.67	0.41
1:A:19:VAL:HG11	1:A:407:TYR:CZ	2.55	0.41
2:D:241:LYS:NZ	2:D:256:ASP:OD2	2.49	0.41
2:D:316:LEU:HD11	2:D:331:LYS:HG2	2.02	0.41
1:A:42:SER:OG	1:A:384:ASN:HB3	2.20	0.41
2:B:175:ILE:CD1	2:B:181:VAL:HG21	2.51	0.41
1:C:382:ALA:HB1	1:C:386:ASP:HB2	2.03	0.41
1:A:136:ASP:CG	1:A:170:LYS:HZ1	2.24	0.41
1:A:42:SER:CB	1:A:388:ASP:OD1	2.69	0.41
2:B:521:LEU:HD22	2:D:447:TYR:CD1	2.55	0.41
1:A:31:HIS:NE2	1:A:47:ILE:HD11	2.36	0.41
2:B:254:LEU:HD22	2:B:281:MET:SD	2.61	0.41
1:C:52:SER:HB3	5:C:508:GOL:C1	2.50	0.41
2:D:414:PRO:O	2:D:417:LYS:HG3	2.21	0.41
1:A:22:GLU:HG2	1:A:26:LYS:HE3	2.02	0.41
1:A:467:LEU:HD23	1:A:467:LEU:C	2.41	0.41
1:C:68:LYS:NZ	5:C:505:GOL:H31	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:422:GLY:HA2	1:C:439:ARG:O	2.21	0.41
1:C:447:SER:OG	1:C:448:GLY:N	2.53	0.41
2:B:113:CYS:CB	9:B:947:HOH:O	2.69	0.41
2:B:296:TRP:CD1	2:B:377:MET:HE3	2.55	0.41
1:C:62:CYS:SG	1:C:64:TYR:HB3	2.61	0.41
2:D:151:THR:HG22	2:D:162:LEU:HD21	2.03	0.41
1:A:274:HIS:CD2	1:A:451:HIS:CE1	3.08	0.41
1:C:22:GLU:OE2	1:C:25:ARG:NH2	2.54	0.41
1:C:39:VAL:HG23	1:C:391:MET:SD	2.61	0.41
1:C:440:GLU:HG3	9:C:609:HOH:O	2.20	0.41
1:C:68:LYS:HZ1	5:C:505:GOL:H31	1.86	0.41
1:A:355:ILE:HB	1:A:360:PRO:HD3	2.02	0.41
2:B:151:THR:CG2	2:B:162:LEU:HD11	2.50	0.41
2:B:48:THR:OG1	2:B:458:LYS:HE3	2.21	0.41
2:D:88:TYR:O	2:D:149:VAL:HA	2.21	0.41
1:A:428:LYS:CB	9:A:785:HOH:O	2.51	0.40
1:C:135:ILE:HD13	1:C:178:ILE:HD13	2.03	0.40
5:B:603:GOL:O2	2:D:514:ALA:O	2.35	0.40
1:A:253:TRP:CZ2	1:A:262:ILE:HG23	2.57	0.40
1:A:277:ARG:HD2	1:A:386:ASP:OD2	2.20	0.40
2:B:375:PHE:HD1	5:B:607:GOL:C3	2.35	0.40
1:C:379:TYR:CG	1:C:382:ALA:HB2	2.57	0.40
2:D:449:LYS:HB2	2:D:449:LYS:HE3	1.79	0.40
1:A:97:ARG:O	1:A:231:ILE:HA	2.22	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	479/490 (98%)	457 (95%)	20 (4%)	2 (0%)	34 22

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	477/490 (97%)	448 (94%)	29 (6%)	0	100	100
2	B	521/522 (100%)	504 (97%)	16 (3%)	1 (0%)	47	38
2	D	520/522 (100%)	501 (96%)	18 (4%)	1 (0%)	47	38
All	All	1997/2024 (99%)	1910 (96%)	83 (4%)	4 (0%)	47	38

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	94	ALA
2	B	255	SER
2	D	255	SER
1	A	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	409/413 (99%)	405 (99%)	4 (1%)	76	73
1	C	408/413 (99%)	402 (98%)	6 (2%)	65	59
2	B	455/454 (100%)	452 (99%)	3 (1%)	84	83
2	D	454/454 (100%)	450 (99%)	4 (1%)	78	77
All	All	1726/1734 (100%)	1709 (99%)	17 (1%)	76	73

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	92	SER
1	A	98	ASN
1	A	362	HIS
1	A	401	TYR
2	B	177	ASP
2	B	258	GLU
2	B	449	LYS

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Mol	Chain	Res	Type
1	C	43	LYS
1	C	98	ASN
1	C	279	MET
1	C	362	HIS
1	C	401	TYR
1	C	445	ASP
2	D	88	TYR
2	D	102	TYR
2	D	172	GLU
2	D	258	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	274	HIS
1	A	384	ASN
2	B	65	ASN
1	C	98	ASN
1	C	321	GLN
1	C	384	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 25 ligands modelled in this entry, 2 are monoatomic - leaving 23 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	$\# Z > 2$
5	GOL	C	504	-	5,5,5	0.43	0	5,5,5	0.56	0
6	1CL	C	509	1,2	0,22,22	0.00	-	-		
5	GOL	D	601	-	5,5,5	0.40	0	5,5,5	0.21	0
5	GOL	B	606	-	5,5,5	0.40	0	5,5,5	0.98	0
5	GOL	B	603	-	5,5,5	0.46	0	5,5,5	0.74	0
5	GOL	D	603	-	5,5,5	0.38	0	5,5,5	0.40	0
6	1CL	A	504	1,2	0,22,22	0.00	-	-		
5	GOL	B	608	-	5,5,5	1.11	1 (20%)	5,5,5	2.28	1 (20%)
3	HCA	A	501	-	4,13,13	1.16	0	4,18,18	1.72	1 (25%)
4	ICS	A	502	1	18,30,30	7.77	13 (72%)	-		
5	GOL	C	506	-	5,5,5	0.49	0	5,5,5	0.54	0
5	GOL	A	503	-	5,5,5	0.18	0	5,5,5	0.63	0
5	GOL	C	508	-	5,5,5	0.63	0	5,5,5	0.71	0
3	HCA	C	501	-	4,13,13	0.93	0	4,18,18	2.52	2 (50%)
5	GOL	C	505	-	5,5,5	0.99	1 (20%)	5,5,5	1.67	1 (20%)
4	ICS	C	502	1	18,30,30	8.20	15 (83%)	-		
8	TRS	B	604	-	7,7,7	0.43	0	9,9,9	1.09	0
5	GOL	C	503	-	5,5,5	0.47	0	5,5,5	1.42	1 (20%)
5	GOL	B	607	-	5,5,5	0.31	0	5,5,5	0.93	0
5	GOL	C	507	-	5,5,5	0.86	0	5,5,5	1.33	0
8	TRS	D	602	-	7,7,7	0.27	0	9,9,9	0.66	0
5	GOL	B	602	-	5,5,5	0.64	0	5,5,5	0.82	0
5	GOL	B	605	-	5,5,5	0.23	0	5,5,5	0.86	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
6	1CL	C	509	1,2	-	-	0/9/8/8
5	GOL	B	606	-	-	2/4/4/4	-
8	TRS	B	604	-	-	3/9/9/9	-
5	GOL	C	506	-	-	2/4/4/4	-
6	1CL	A	504	1,2	-	-	0/9/8/8

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GOL	B	603	-	-	2/4/4/4	-
5	GOL	C	503	-	-	3/4/4/4	-
5	GOL	B	607	-	-	2/4/4/4	-
5	GOL	D	603	-	-	2/4/4/4	-
5	GOL	C	504	-	-	4/4/4/4	-
8	TRS	D	602	-	-	7/9/9/9	-
3	HCA	A	501	-	-	4/7/17/17	-
5	GOL	C	508	-	-	2/4/4/4	-
3	HCA	C	501	-	-	2/7/17/17	-
5	GOL	B	608	-	-	4/4/4/4	-
5	GOL	C	505	-	-	2/4/4/4	-
5	GOL	A	503	-	-	4/4/4/4	-
5	GOL	C	507	-	-	2/4/4/4	-
5	GOL	D	601	-	-	1/4/4/4	-
5	GOL	B	602	-	-	2/4/4/4	-
5	GOL	B	605	-	-	4/4/4/4	-

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	502	ICS	S4A-FE3	-11.74	2.03	2.32
4	A	502	ICS	S4A-FE3	-10.90	2.05	2.32
4	C	502	ICS	S4B-FE7	-10.89	2.05	2.32
4	C	502	ICS	S1B-FE6	-10.83	2.05	2.32
4	A	502	ICS	S1A-FE4	-10.48	2.06	2.32
4	C	502	ICS	S4A-FE4	-10.36	2.07	2.32
4	A	502	ICS	S3B-FE7	-10.35	2.07	2.32
4	C	502	ICS	S4B-FE5	-10.23	2.07	2.32
4	C	502	ICS	S1A-FE4	-9.76	2.08	2.32
4	A	502	ICS	S3B-FE6	-9.66	2.08	2.32
4	C	502	ICS	S3B-FE6	-9.60	2.08	2.32
4	C	502	ICS	S1B-FE5	-9.54	2.09	2.32
4	C	502	ICS	S1A-FE2	-9.42	2.09	2.32
4	C	502	ICS	S3B-FE7	-9.39	2.09	2.32
4	A	502	ICS	S2A-FE3	-9.36	2.09	2.32
4	A	502	ICS	S1B-FE6	-9.31	2.09	2.32
4	A	502	ICS	S4B-FE7	-9.26	2.09	2.32
4	A	502	ICS	S1A-FE2	-9.13	2.10	2.32
4	A	502	ICS	S4B-FE5	-9.07	2.10	2.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	502	ICS	S1B-FE5	-8.85	2.10	2.32
4	A	502	ICS	S4A-FE4	-8.56	2.11	2.32
4	C	502	ICS	S2A-FE2	-8.55	2.11	2.32
4	A	502	ICS	S2A-FE2	-8.01	2.12	2.32
4	C	502	ICS	S2A-FE3	-7.82	2.13	2.32
4	C	502	ICS	S5A-FE3	-3.95	2.15	2.24
4	C	502	ICS	S3A-FE4	-3.56	2.16	2.24
4	A	502	ICS	S5A-FE3	-3.38	2.16	2.24
5	B	608	GOL	O3-C3	2.48	1.52	1.42
4	C	502	ICS	S3A-FE5	-2.10	2.19	2.24
5	C	505	GOL	O2-C2	2.01	1.49	1.43

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	501	HCA	C4-C5-C6	4.07	117.55	111.39
5	B	608	GOL	C3-C2-C1	3.68	126.00	111.70
3	C	501	HCA	C3-C2-C1	-2.84	110.43	114.98
3	A	501	HCA	C3-C2-C1	-2.38	111.17	114.98
5	C	505	GOL	O3-C3-C2	2.26	121.04	110.20
5	C	503	GOL	O2-C2-C3	2.11	118.42	109.12

There are no chirality outliers.

All (54) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	504	GOL	O1-C1-C2-C3
5	C	504	GOL	C1-C2-C3-O3
5	B	603	GOL	O1-C1-C2-C3
5	B	608	GOL	O2-C2-C3-O3
3	A	501	HCA	C2-C3-C4-C5
3	A	501	HCA	C7-C3-C4-C5
5	A	503	GOL	O1-C1-C2-C3
5	A	503	GOL	C1-C2-C3-O3
5	C	508	GOL	C1-C2-C3-O3
5	C	503	GOL	C1-C2-C3-O3
8	D	602	TRS	C2-C-C1-O1
8	D	602	TRS	N-C-C1-O1
5	B	605	GOL	C1-C2-C3-O3
5	B	603	GOL	O1-C1-C2-O2
5	B	608	GOL	O1-C1-C2-O2
5	C	506	GOL	O2-C2-C3-O3

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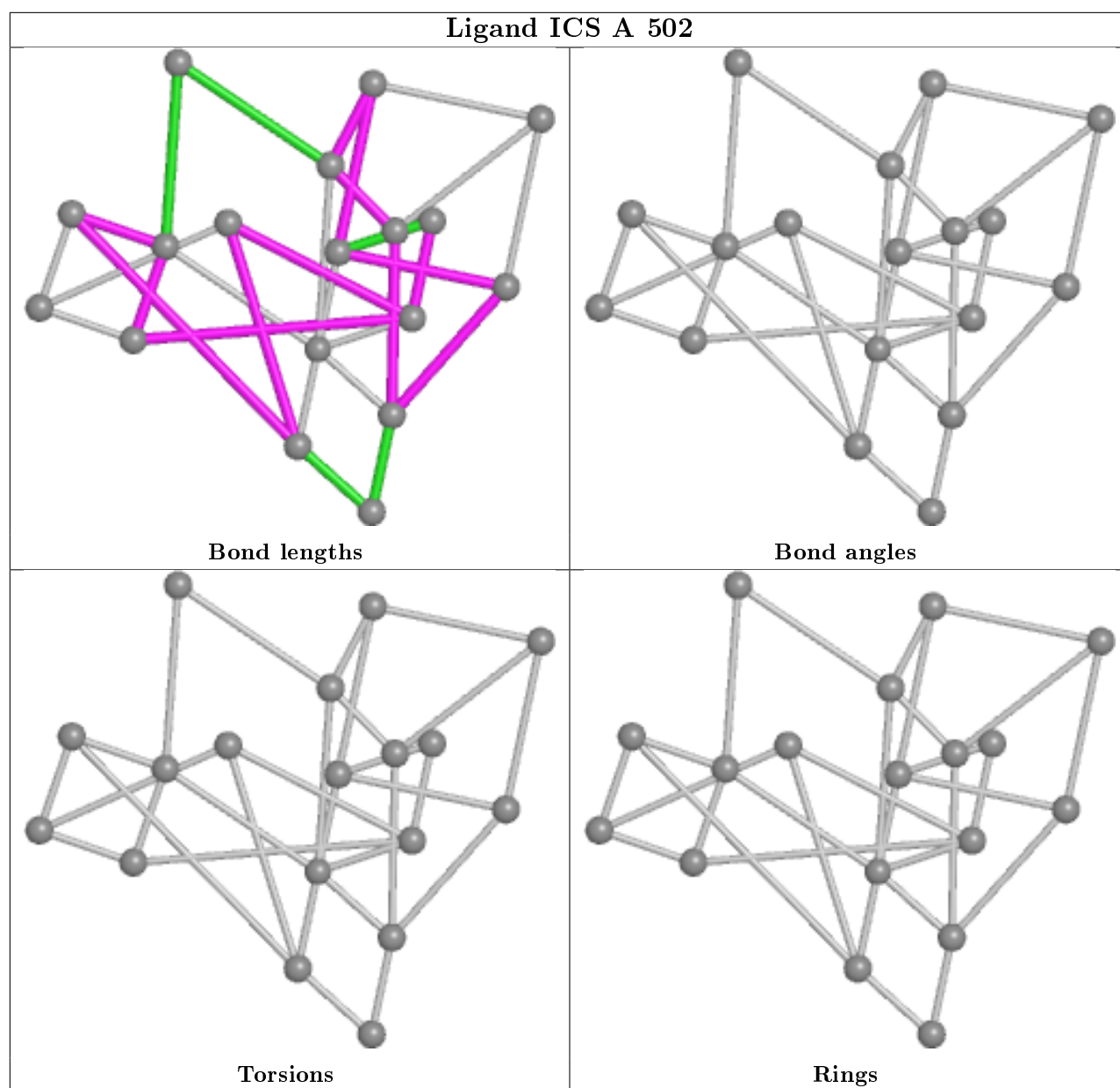
Mol	Chain	Res	Type	Atoms
5	A	503	GOL	O1-C1-C2-O2
5	A	503	GOL	O2-C2-C3-O3
5	B	606	GOL	O1-C1-C2-C3
5	D	603	GOL	O1-C1-C2-C3
5	B	608	GOL	O1-C1-C2-C3
5	B	608	GOL	C1-C2-C3-O3
5	C	506	GOL	C1-C2-C3-O3
5	C	505	GOL	C1-C2-C3-O3
5	C	503	GOL	O1-C1-C2-C3
5	B	607	GOL	C1-C2-C3-O3
5	C	507	GOL	C1-C2-C3-O3
5	B	602	GOL	C1-C2-C3-O3
5	B	605	GOL	O1-C1-C2-C3
5	C	504	GOL	O1-C1-C2-O2
5	C	504	GOL	O2-C2-C3-O3
5	C	508	GOL	O2-C2-C3-O3
5	B	607	GOL	O2-C2-C3-O3
5	B	605	GOL	O2-C2-C3-O3
8	D	602	TRS	C3-C-C1-O1
5	C	503	GOL	O2-C2-C3-O3
5	B	602	GOL	O2-C2-C3-O3
5	C	507	GOL	O2-C2-C3-O3
5	B	605	GOL	O1-C1-C2-O2
8	B	604	TRS	C1-C-C3-O3
8	B	604	TRS	C2-C-C3-O3
8	D	602	TRS	C2-C-C3-O3
5	D	603	GOL	O1-C1-C2-O2
3	A	501	HCA	C1-C2-C3-C4
8	D	602	TRS	C1-C-C3-O3
5	D	601	GOL	O2-C2-C3-O3
5	B	606	GOL	O1-C1-C2-O2
3	C	501	HCA	C1-C2-C3-C4
5	C	505	GOL	O1-C1-C2-O2
3	A	501	HCA	O7-C3-C4-C5
8	B	604	TRS	N-C-C3-O3
8	D	602	TRS	C1-C-C2-O2
8	D	602	TRS	N-C-C3-O3
3	C	501	HCA	C2-C3-C4-C5

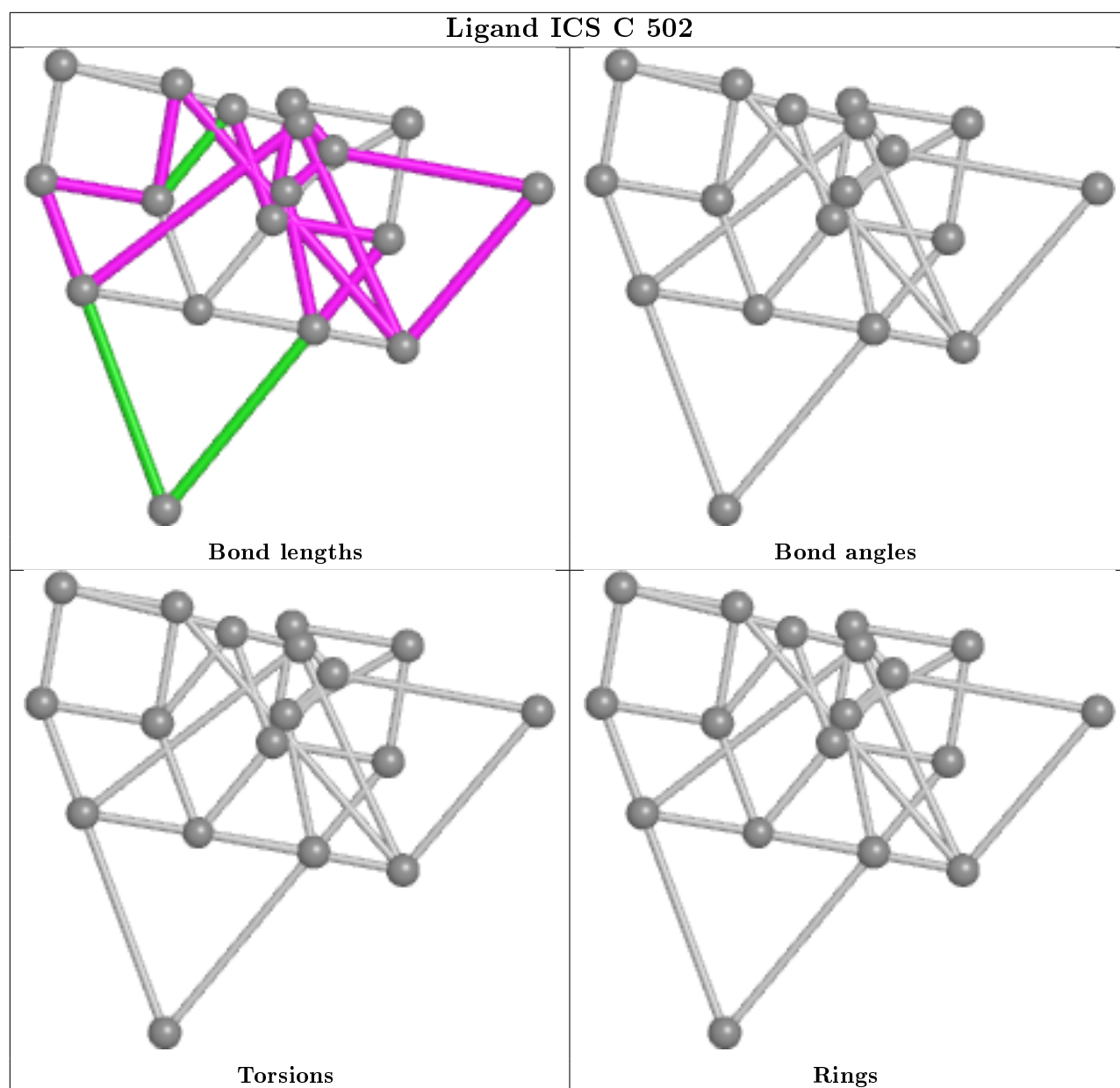
There are no ring outliers.

14 monomers are involved in 42 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	504	GOL	3	0
5	B	606	GOL	1	0
5	B	603	GOL	3	0
5	B	608	GOL	7	0
3	A	501	HCA	1	0
5	C	508	GOL	10	0
3	C	501	HCA	2	0
5	C	505	GOL	4	0
8	B	604	TRS	2	0
5	B	607	GOL	4	0
5	C	507	GOL	2	0
8	D	602	TRS	1	0
5	B	602	GOL	1	0
5	B	605	GOL	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.





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, 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	479/490 (97%)	0.87	56 (11%) 4 5	24, 39, 65, 100	12 (2%)
1	C	478/490 (97%)	1.11	78 (16%) 1 1	28, 46, 76, 93	4 (0%)
2	B	522/522 (100%)	0.46	22 (4%) 36 38	24, 36, 53, 70	5 (0%)
2	D	522/522 (100%)	0.44	21 (4%) 38 40	22, 37, 55, 80	13 (2%)
All	All	2001/2024 (98%)	0.71	177 (8%) 10 11	22, 39, 64, 100	34 (1%)

All (177) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	123	ALA	8.1
2	B	123	ALA	5.9
2	D	121	ASP	5.5
1	C	39	VAL	5.4
1	A	3	GLY	5.0
1	C	14	GLN	5.0
1	A	481	ALA	4.5
1	C	172	ALA	4.4
1	C	318	GLU	4.2
1	A	39	VAL	4.2
1	C	217	ALA	4.1
1	C	168	LYS	4.0
1	C	36	ASP	4.0
2	B	121	ASP	3.9
1	A	92	SER	3.9
1	C	7	GLU	3.7
1	A	124	VAL	3.6
2	B	101	SER	3.6
1	A	18	GLU	3.6
1	C	392	LYS	3.6
1	C	92	SER	3.6

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Mol	Chain	Res	Type	RSRZ
1	C	125	PHE	3.6
2	B	178	GLU	3.6
1	C	443[A]	SER	3.5
1	C	322	LYS	3.5
1	C	317	ASP	3.5
1	C	44	LYS	3.4
1	C	59	ILE	3.4
2	B	177	ASP	3.4
1	C	292	ILE	3.4
1	C	75	ILE	3.4
2	D	215	ASP	3.3
1	C	221	TYR	3.3
1	A	212	GLU	3.3
1	A	125	PHE	3.2
2	B	312	GLU	3.2
1	A	75	ILE	3.2
1	A	40	THR	3.2
1	A	59	ILE	3.2
1	C	163	ILE	3.2
2	D	125	PHE	3.2
2	B	215	ASP	3.2
2	B	214	ASP	3.2
1	C	70	VAL	3.2
2	B	521	LEU	3.2
1	A	43	LYS	3.2
1	C	18	GLU	3.1
1	C	231	ILE	3.1
1	A	231	ILE	3.1
1	C	257	GLY	3.1
1	A	322	LYS	3.1
1	A	19	VAL	3.1
1	A	391	MET	3.1
1	A	70	VAL	3.1
1	A	45	CYS	3.1
1	A	392	LYS	3.1
1	C	24	ALA	3.1
1	C	291	GLY	3.1
1	C	15	GLU	3.0
2	B	96	VAL	3.0
2	D	417	LYS	3.0
1	A	409	PHE	3.0
2	D	101	SER	3.0

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Mol	Chain	Res	Type	RSRZ
2	B	417	LYS	3.0
1	A	44	LYS	3.0
2	D	96	VAL	3.0
1	C	194	GLY	2.9
2	B	124	VAL	2.9
1	C	391	MET	2.9
1	A	443[A]	SER	2.9
1	C	162	ASP	2.9
1	C	43	LYS	2.8
1	A	168	LYS	2.8
1	C	40	THR	2.8
2	B	173	GLY	2.8
2	D	521	LEU	2.8
1	C	425	ILE	2.8
1	A	215	THR	2.7
1	A	279	MET	2.7
1	A	7	GLU	2.7
1	A	38	ALA	2.7
2	D	217	VAL	2.7
1	A	174	LEU	2.7
1	C	320	ILE	2.7
2	B	125	PHE	2.7
1	C	319	SER	2.7
2	D	124	VAL	2.7
1	C	176	LYS	2.7
1	C	212	GLU	2.7
1	C	94	ALA	2.6
1	A	318	GLU	2.6
2	D	102	TYR	2.6
1	A	172	ALA	2.6
1	A	41	GLN	2.6
1	C	476	GLN	2.6
2	B	38	ASP	2.6
2	D	38	ASP	2.6
1	C	321	GLN	2.6
1	A	72	TRP	2.5
1	C	157	GLY	2.5
1	A	291	GLY	2.5
1	A	42	SER	2.5
1	C	161	ASP	2.5
2	D	450	PHE	2.5
1	C	200	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	259	ILE	2.5
1	C	110	VAL	2.5
1	C	400	LEU	2.5
1	C	409	PHE	2.5
2	B	447	TYR	2.5
1	C	46	ILE	2.4
1	C	288	GLU	2.4
2	B	168	ASN	2.4
2	D	89	VAL	2.4
2	D	157	VAL	2.4
1	C	69	GLY	2.4
1	C	98	ASN	2.4
1	C	41	GLN	2.4
1	C	366	ALA	2.4
2	D	214	ASP	2.4
1	A	126	GLY	2.4
1	C	316	PHE	2.3
1	C	3	GLY	2.3
1	A	8	GLU	2.3
1	C	475	LEU	2.3
1	A	93	ARG	2.3
1	C	175	SER	2.3
1	C	37	PRO	2.3
2	B	450	PHE	2.2
1	C	123	ILE	2.2
2	D	46	TRP	2.2
1	A	257	GLY	2.2
1	A	36	ASP	2.2
1	C	16	VAL	2.2
1	A	477	ALA	2.2
2	D	265	ALA	2.2
2	B	261	LEU	2.2
1	A	15	GLU	2.2
1	C	190	SER	2.2
1	A	110	VAL	2.2
1	C	174	LEU	2.2
1	A	407	TYR	2.2
1	C	356	GLY	2.2
1	A	449	PRO	2.1
1	C	397	SER	2.1
1	C	327	VAL	2.1
1	C	56	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	71	VAL	2.1
1	C	314	ALA	2.1
1	C	477	ALA	2.1
2	D	261	LEU	2.1
1	C	93	ARG	2.1
1	A	213	ASP	2.1
1	C	253	TRP	2.1
1	A	4	MET	2.1
1	A	9	VAL	2.1
1	C	4	MET	2.1
1	A	5	SER	2.1
1	A	175	SER	2.1
1	C	45	CYS	2.1
2	B	153	CYS	2.1
2	D	122	ALA	2.1
1	A	157	GLY	2.1
2	B	102	TYR	2.1
1	C	72	TRP	2.1
1	A	11	SER	2.1
1	A	217	ALA	2.1
1	C	225	ILE	2.1
1	A	112	MET	2.0
1	C	128	ASP	2.0
1	A	397	SER	2.0
2	D	168	ASN	2.0
1	A	94	ALA	2.0
1	C	34	VAL	2.0
2	B	77	LEU	2.0
1	A	173	GLU	2.0
1	C	480	GLU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

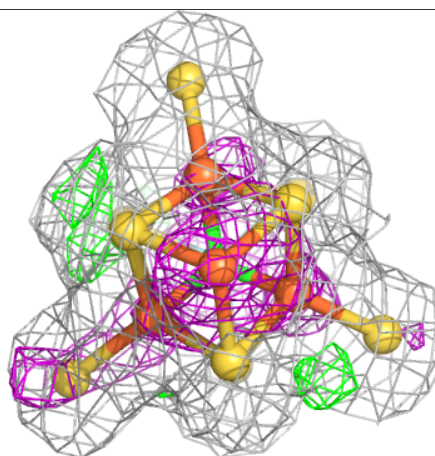
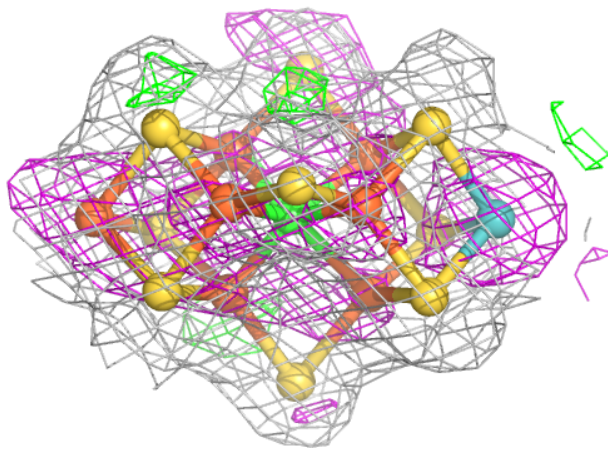
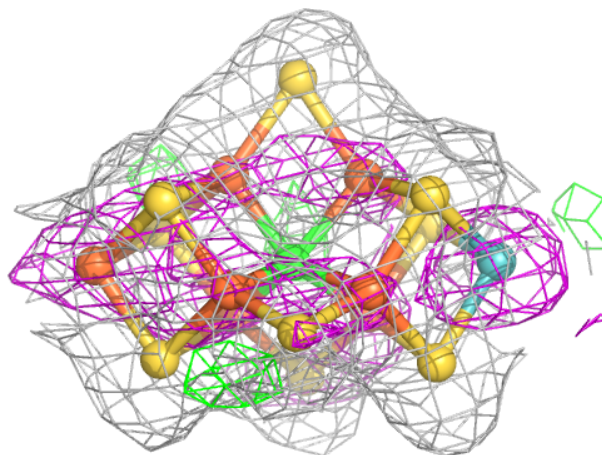
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	GOL	C	508	6/6	0.62	0.44	54,63,70,81	0
8	TRS	B	604	8/8	0.64	0.25	53,60,71,72	0
5	GOL	C	505	6/6	0.70	0.35	34,43,48,52	0
5	GOL	C	503	6/6	0.71	0.20	48,58,61,63	0
5	GOL	B	602	6/6	0.71	0.26	40,48,53,55	0
5	GOL	A	503	6/6	0.72	0.18	39,42,44,44	0
5	GOL	C	507	6/6	0.73	0.20	49,52,57,59	0
8	TRS	D	602	8/8	0.74	0.20	52,60,65,66	0
5	GOL	C	504	6/6	0.76	0.20	42,51,56,58	0
5	GOL	D	601	6/6	0.76	0.17	54,57,59,65	0
5	GOL	B	603	6/6	0.80	0.15	37,38,40,46	0
5	GOL	B	608	6/6	0.83	0.33	24,35,53,63	0
5	GOL	C	506	6/6	0.84	0.33	57,72,80,83	0
5	GOL	B	606	6/6	0.86	0.22	47,53,57,69	0
5	GOL	D	603	6/6	0.86	0.16	39,48,53,54	0
5	GOL	B	605	6/6	0.90	0.12	37,46,53,53	0
3	HCA	A	501	14/14	0.91	0.15	27,30,39,39	0
3	HCA	C	501	14/14	0.92	0.15	28,34,39,39	0
5	GOL	B	607	6/6	0.92	0.32	35,45,51,52	0
6	1CL	C	509	15/15	0.93	0.08	28,31,40,42	0
6	1CL	A	504	15/15	0.93	0.08	26,30,37,41	15
7	CA	B	609	1/1	0.94	0.04	42,42,42,42	0
4	ICS	C	502	18/18	0.98	0.07	28,33,39,40	0
4	ICS	A	502	18/18	0.98	0.06	25,32,35,37	0
7	CA	B	601	1/1	0.99	0.04	41,41,41,41	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.

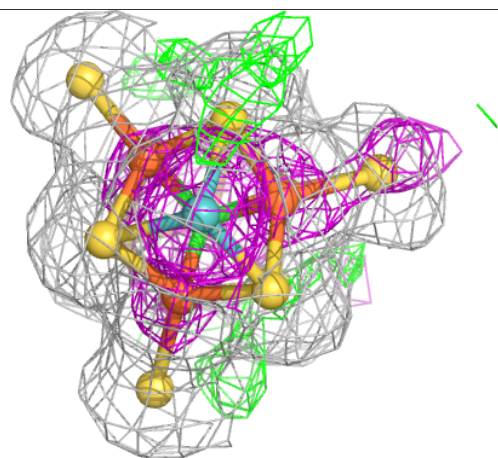
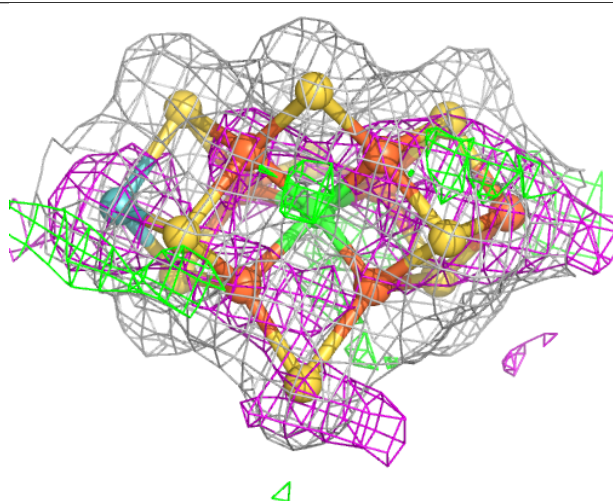
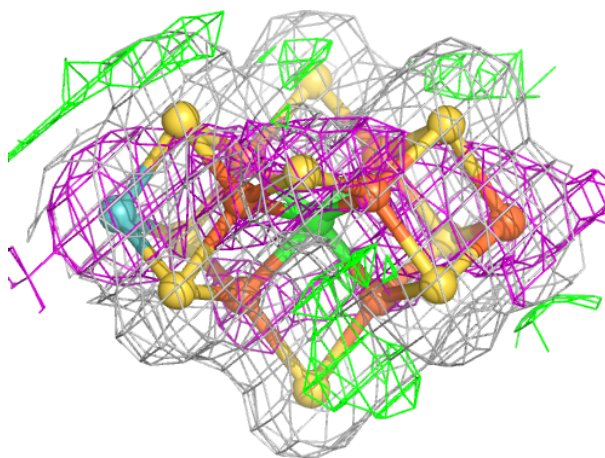
Electron density around ICS C 502:

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



Electron density around ICS A 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 ⓘ

There are no such residues in this entry.