



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 01:29 PM EDT

PDB ID : 4MK4 / pdb_00004mk4
Title : S197C variant of human ferrochelatase.
Authors : Lanzilotta, W.N.; Medlock, A.E.; Dailey, T.E.; Dailey, H.A.
Deposited on : 2013-09-04
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
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

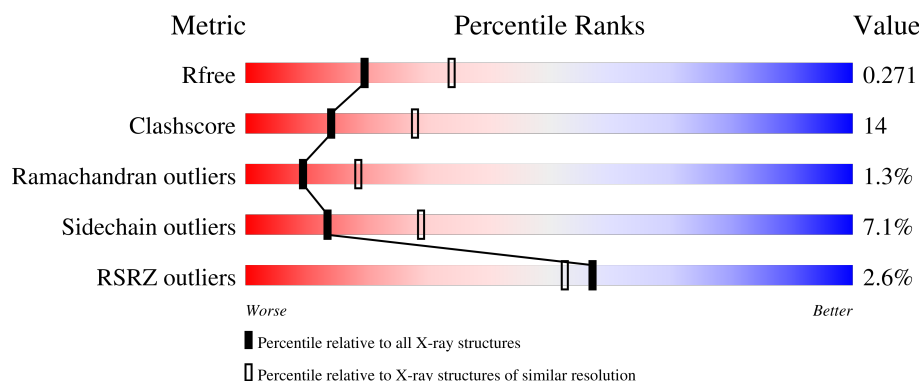
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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

Mol	Chain	Length	Quality of chain
1	A	359	2% 69% 25% 6%
1	B	359	3% 71% 25% .

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6092 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.

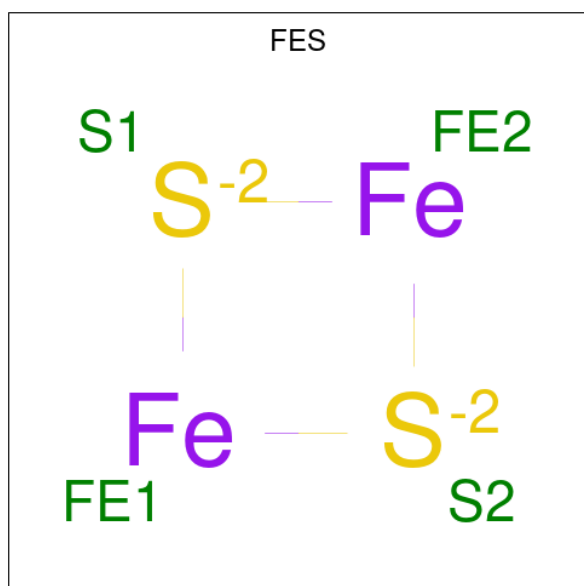
- Molecule 1 is a protein called Ferrochelatase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	359	Total	C	N	O	S	0	0	0
			2891	1841	503	528	19			
1	B	359	Total	C	N	O	S	0	1	0
			2900	1846	505	530	19			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	197	CYS	SER	engineered mutation	UNP P22830
B	197	CYS	SER	engineered mutation	UNP P22830

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂).



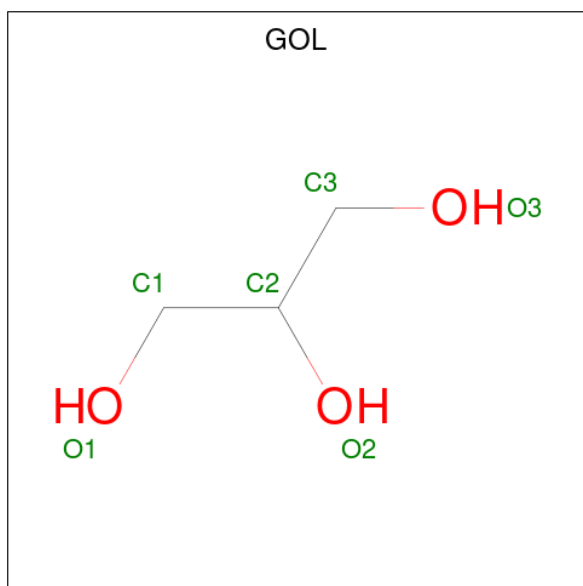
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			4	2	2		

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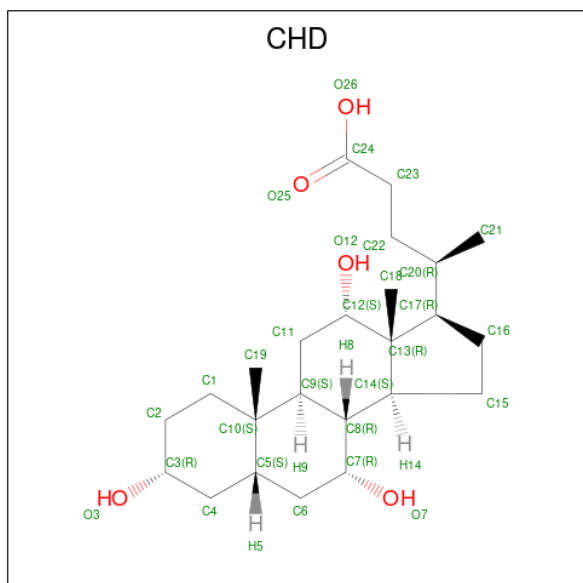
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	Fe	S		
2	B	1	4	2	2	0	0

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 4 is CHOLIC ACID (CCD ID: CHD) (formula: $C_{24}H_{40}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			29	24	5		
4	A	1	Total	C	O	0	0
			29	24	5		
4	B	1	Total	C	O	0	0
			29	24	5		
4	B	1	Total	C	O	0	0
			29	24	5		
4	B	1	Total	C	O	0	0
			29	24	5		

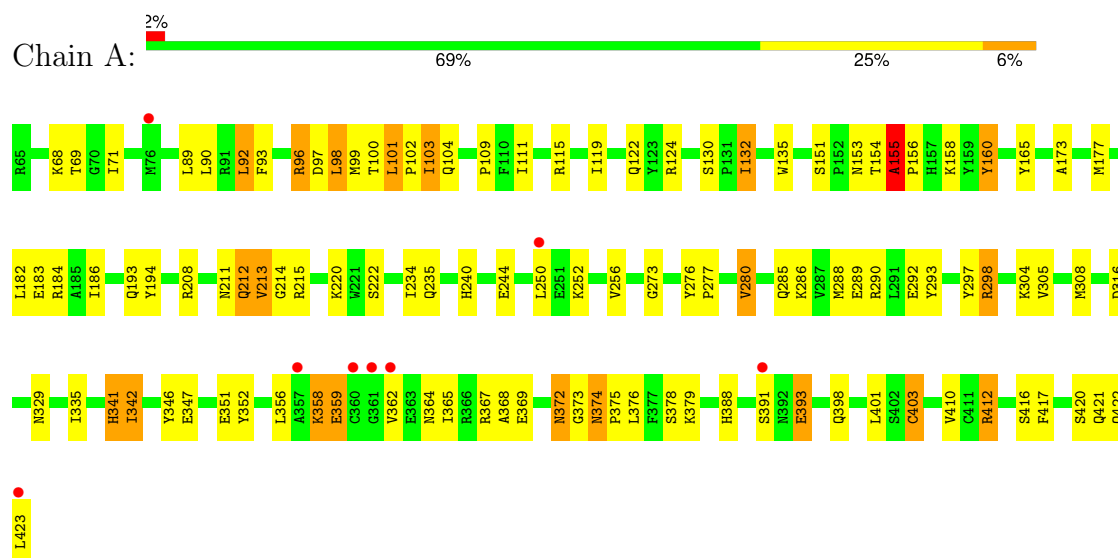
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	79	Total	O	0	0
			79	79		
5	B	63	Total	O	0	0
			63	63		

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: Ferrochelatase, mitochondrial



• Molecule 1: Ferrochelatase, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.17Å 93.43Å 111.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.00 – 2.50 60.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	94.2 (60.00-2.50) 94.2 (60.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.29 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.221 , 0.277 0.217 , 0.271	Depositor DCC
R_{free} test set	1563 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	35.2	Xtriage
Anisotropy	0.522	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6092	wwPDB-VP
Average B, all atoms (Å ²)	37.0	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 67.14 % 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 5.3960e-06. 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: FES, GOL, CHD

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	0.91	0/2961	1.13	7/4011 (0.2%)
1	B	0.91	1/2970 (0.0%)	1.10	5/4023 (0.1%)
All	All	0.91	1/5931 (0.0%)	1.12	12/8034 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	362	VAL	CA-CB	6.66	1.63	1.54

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	410	VAL	N-CA-C	6.60	117.38	110.72
1	A	280	VAL	N-CA-C	-6.15	104.35	110.62
1	A	155	ALA	CA-C-N	-5.92	113.44	120.13
1	A	155	ALA	C-N-CA	-5.92	113.44	120.13
1	A	92	LEU	CA-CB-CG	5.74	136.40	116.30
1	B	364	ASN	N-CA-C	5.53	122.58	110.80
1	B	194	TYR	N-CA-C	5.29	117.36	109.59
1	B	377	PHE	N-CA-C	-5.22	105.65	112.23
1	A	186	ILE	N-CA-C	5.15	115.00	107.37
1	B	420	SER	N-CA-C	5.12	119.51	113.16
1	B	280	VAL	N-CA-C	-5.06	104.85	110.62
1	A	365	ILE	CB-CA-C	5.04	117.23	110.42

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	2891	0	2899	85	0
1	B	2900	0	2906	79	0
2	A	4	0	0	0	0
2	B	4	0	0	1	0
3	A	6	0	8	0	0
4	A	58	0	78	6	0
4	B	87	0	117	13	0
5	A	79	0	0	4	0
5	B	63	0	0	4	0
All	All	6092	0	6008	171	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 14.

All (171) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:304:LYS:HE2	5:A:649:HOH:O	1.44	1.13
1:B:300:VAL:HG12	1:B:313:PRO:HG2	1.37	1.05
1:A:155:ALA:HB1	1:A:156:PRO:HD3	1.32	1.03
1:B:235:GLN:HG3	1:B:290:ARG:NH1	1.79	0.98
1:B:155:ALA:HB1	1:B:156:PRO:CD	1.93	0.97
1:A:155:ALA:HB1	1:A:156:PRO:CD	1.96	0.96
1:B:155:ALA:HB1	1:B:156:PRO:HD3	1.49	0.92
1:A:374:ASN:ND2	1:A:376:LEU:H	1.67	0.92
1:B:240:HIS:HD2	1:B:369:GLU:O	1.55	0.88
1:B:244:GLU:HG3	1:B:368:ALA:HA	1.56	0.85
1:A:222:SER:OG	1:A:388:HIS:HE1	1.58	0.84
1:A:398:GLN:HE22	1:B:297:TYR:H	1.27	0.81
1:A:240:HIS:HD2	1:A:369:GLU:O	1.62	0.81
1:A:115:ARG:HG2	1:A:115:ARG:HH11	1.48	0.79
1:A:297:TYR:H	1:B:398:GLN:HE22	1.28	0.77
1:B:222:SER:OG	1:B:388:HIS:HE1	1.68	0.77
1:A:374:ASN:HD22	1:A:376:LEU:H	1.32	0.77
1:A:244:GLU:HG3	1:A:368:ALA:HA	1.67	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:ALA:CB	1:A:156:PRO:CD	2.64	0.76
1:A:288:MET:HG3	1:A:297:TYR:CD2	2.21	0.75
1:A:374:ASN:HD22	1:A:374:ASN:C	1.94	0.75
1:B:68:LYS:CB	1:B:155:ALA:HB3	2.18	0.73
1:B:68:LYS:HB3	1:B:155:ALA:HB3	1.69	0.72
1:A:222:SER:OG	1:A:388:HIS:CE1	2.44	0.71
1:B:155:ALA:CB	1:B:156:PRO:CD	2.69	0.69
1:B:123:TYR:CZ	1:B:342:ILE:HG13	2.27	0.69
1:B:319:ILE:HD11	1:B:332:LEU:HD21	1.76	0.68
1:B:240:HIS:CD2	1:B:369:GLU:O	2.44	0.67
1:B:244:GLU:HG2	5:B:655:HOH:O	1.95	0.66
1:B:328:LYS:O	1:B:363:GLU:O	2.13	0.66
4:A:504:CHD:H12	4:A:504:CHD:H212	1.79	0.65
1:B:114:ARG:HG2	1:B:114:ARG:HH11	1.60	0.64
1:B:68:LYS:HD3	1:B:155:ALA:CB	2.27	0.64
1:A:329:ASN:HD22	1:A:364:ASN:HB2	1.63	0.64
1:A:102:PRO:O	1:A:103:ILE:HG12	1.98	0.63
1:A:250:LEU:HD12	1:A:250:LEU:H	1.64	0.62
1:B:171:GLU:CD	1:B:171:GLU:H	2.07	0.62
1:B:330:ILE:HD12	1:B:330:ILE:N	2.14	0.62
1:A:234:ILE:HG13	1:A:286:LYS:HG2	1.81	0.62
1:A:173:ALA:O	1:A:177:MET:HG3	2.00	0.62
1:A:89:LEU:HA	1:A:92:LEU:HD23	1.82	0.62
1:B:222:SER:OG	1:B:388:HIS:CE1	2.51	0.62
1:A:89:LEU:HA	1:A:92:LEU:CD2	2.29	0.62
1:A:358:LYS:O	1:A:359:GLU:CB	2.48	0.61
1:B:115:ARG:HD2	4:B:502:CHD:H221	1.83	0.61
1:B:235:GLN:HG3	1:B:290:ARG:HH11	1.63	0.60
1:A:69:THR:HG23	1:A:184:ARG:HD2	1.82	0.60
1:A:412:ARG:HB2	1:A:412:ARG:HH21	1.66	0.60
1:A:417:PHE:O	1:A:421:GLN:HG2	2.03	0.59
1:A:98:LEU:HD11	1:A:165:TYR:OH	2.03	0.58
1:B:68:LYS:HD3	1:B:155:ALA:HB2	1.84	0.58
1:A:297:TYR:C	1:A:297:TYR:CD1	2.82	0.58
1:A:374:ASN:HD22	1:A:376:LEU:N	2.00	0.58
1:A:374:ASN:HD22	1:A:375:PRO:N	2.01	0.57
1:B:416:SER:O	1:B:420:SER:HB2	2.04	0.57
1:B:422:GLN:O	1:B:423:LEU:HB2	2.04	0.57
1:B:374:ASN:HD22	1:B:376:LEU:H	1.53	0.57
1:B:288:MET:HG3	1:B:297:TYR:CD2	2.40	0.56
1:A:93:PHE:CE1	1:A:111:ILE:HG22	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:374:ASN:ND2	1:A:376:LEU:N	2.46	0.56
1:B:111:ILE:HG23	4:B:502:CHD:H213	1.89	0.55
1:B:374:ASN:ND2	1:B:376:LEU:H	2.04	0.55
1:A:401:LEU:O	1:B:298:ARG:NH2	2.40	0.54
1:B:272:ARG:NH1	2:B:501:FES:S2	2.81	0.54
4:A:503:CHD:H222	5:A:613:HOH:O	2.07	0.54
1:B:354[B]:GLN:O	1:B:358:LYS:HE3	2.07	0.54
1:B:388:HIS:HD2	5:B:656:HOH:O	1.91	0.54
1:B:93:PHE:CE1	1:B:111:ILE:HG22	2.42	0.54
4:A:503:CHD:C22	5:A:613:HOH:O	2.54	0.54
1:A:115:ARG:HG2	1:A:115:ARG:NH1	2.17	0.53
1:A:97:ASP:OD1	1:A:208:ARG:NH2	2.41	0.53
1:A:288:MET:HG3	1:A:297:TYR:CE2	2.42	0.53
1:B:90:LEU:HD13	1:B:109:PRO:HA	1.89	0.53
1:B:89:LEU:HD21	1:B:119:ILE:HD13	1.89	0.52
1:A:93:PHE:HD2	1:A:99:MET:HE1	1.75	0.52
1:A:71:ILE:N	1:A:71:ILE:HD12	2.25	0.51
1:A:194:TYR:HA	5:A:636:HOH:O	2.09	0.51
1:A:235:GLN:HG3	1:A:290:ARG:NH1	2.26	0.51
1:A:101:LEU:HB2	1:A:104:GLN:HG3	1.92	0.50
1:B:263:HIS:HD2	1:B:264:SER:O	1.93	0.50
1:B:114:ARG:HG2	1:B:114:ARG:NH1	2.26	0.50
1:A:391:SER:OG	1:A:393:GLU:HB2	2.11	0.50
1:A:122:GLN:HG2	1:A:346:TYR:CE2	2.47	0.49
1:B:288:MET:HG3	1:B:297:TYR:CE2	2.47	0.49
1:A:358:LYS:O	1:A:359:GLU:HB3	2.12	0.49
1:A:403:CYS:HB3	1:B:298:ARG:NH1	2.27	0.49
1:A:68:LYS:CB	1:A:155:ALA:HB3	2.42	0.49
1:B:68:LYS:HD3	1:B:155:ALA:HB3	1.94	0.49
4:A:504:CHD:H212	4:A:504:CHD:C12	2.42	0.49
1:B:371:LEU:O	1:B:372:ASN:C	2.56	0.49
1:A:135:TRP:CZ3	1:A:372:ASN:HB3	2.48	0.49
1:A:68:LYS:HB2	1:A:155:ALA:HB3	1.96	0.48
1:A:97:ASP:CG	1:A:208:ARG:HH22	2.21	0.48
1:A:90:LEU:HD13	1:A:109:PRO:HA	1.95	0.48
1:A:416:SER:O	1:A:420:SER:HB2	2.14	0.48
1:B:352:TYR:HA	1:B:356:LEU:HB3	1.94	0.48
1:B:102:PRO:C	1:B:103:ILE:HG13	2.37	0.48
1:B:329:ASN:HD22	1:B:364:ASN:HB2	1.77	0.48
1:A:273:GLY:O	1:B:298:ARG:NH1	2.43	0.48
1:B:111:ILE:CG2	4:B:502:CHD:H213	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:374:ASN:HD22	1:B:374:ASN:C	2.23	0.47
1:B:163:PHE:O	1:B:169:LEU:HA	2.15	0.47
1:A:96:ARG:HH12	1:A:104:GLN:HB3	1.78	0.47
1:A:285:GLN:HG2	1:B:282:ALA:HB1	1.96	0.47
4:B:502:CHD:H193	4:B:504:CHD:H162	1.96	0.47
1:A:316:ASP:HB3	1:A:352:TYR:CE1	2.50	0.46
1:A:215:ARG:HE	1:A:215:ARG:HB2	1.54	0.46
1:B:98:LEU:HD11	1:B:165:TYR:OH	2.16	0.46
1:B:68:LYS:HB2	1:B:155:ALA:HB3	1.97	0.46
1:B:356:LEU:O	1:B:357:ALA:HB3	2.16	0.46
1:B:193:GLN:HG2	1:B:280:VAL:HA	1.97	0.46
1:A:288:MET:HE2	1:A:297:TYR:HD2	1.82	0.45
1:B:130:SER:OG	1:B:132:ILE:HG22	2.16	0.45
1:A:89:LEU:HD21	1:A:119:ILE:CD1	2.46	0.45
1:B:171:GLU:CD	1:B:171:GLU:N	2.74	0.45
1:B:347:GLU:O	1:B:351:GLU:HB2	2.16	0.45
1:B:354[A]:GLN:O	1:B:358:LYS:HE3	2.16	0.45
1:A:193:GLN:HG2	1:A:280:VAL:HA	1.98	0.45
4:B:504:CHD:H162	4:B:504:CHD:H222	1.58	0.45
1:A:289:GLU:HG3	1:A:293:TYR:OH	2.17	0.45
1:A:342:ILE:HD13	1:A:346:TYR:HB3	1.98	0.45
1:A:412:ARG:HH21	1:A:412:ARG:CB	2.30	0.45
1:A:151:SER:OG	1:A:154:THR:OG1	2.30	0.44
1:B:374:ASN:HD22	1:B:375:PRO:N	2.15	0.44
4:B:502:CHD:H211	4:B:502:CHD:H232	1.73	0.44
1:B:221:TRP:HB2	1:B:421:GLN:NE2	2.33	0.44
1:B:137:SER:O	1:B:141:GLU:HG3	2.18	0.44
1:A:256:VAL:HA	1:A:329:ASN:O	2.18	0.44
1:A:347:GLU:O	1:A:351:GLU:HB2	2.17	0.44
1:B:131:PRO:O	1:B:132:ILE:C	2.59	0.44
1:A:240:HIS:CD2	1:A:369:GLU:O	2.54	0.44
1:A:335:ILE:O	1:A:335:ILE:HG13	2.18	0.44
1:B:322:LEU:HB3	1:B:327:ARG:HB2	2.00	0.44
4:A:503:CHD:H212	4:A:503:CHD:H12	2.00	0.43
1:A:252:LYS:HD3	1:A:329:ASN:OD1	2.19	0.43
1:A:102:PRO:C	1:A:103:ILE:HG12	2.43	0.43
1:B:68:LYS:HE2	1:B:154:THR:HA	2.00	0.43
1:B:176:GLU:O	1:B:180:ASP:OD1	2.36	0.43
1:A:297:TYR:HD1	1:A:298:ARG:N	2.17	0.42
1:A:102:PRO:O	1:A:103:ILE:C	2.63	0.42
1:A:130:SER:OG	1:A:132:ILE:HG22	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:ILE:HG23	4:B:502:CHD:C21	2.50	0.42
4:B:503:CHD:H222	4:B:503:CHD:H162	1.72	0.42
1:A:160:TYR:N	1:A:160:TYR:CD1	2.87	0.42
1:A:341:HIS:CD2	1:A:341:HIS:N	2.86	0.42
1:B:253:ARG:NH2	5:B:660:HOH:O	2.44	0.42
1:A:96:ARG:HA	1:A:99:MET:O	2.19	0.42
1:B:230:HIS:CE1	1:B:232:LEU:HB2	2.55	0.42
1:B:286:LYS:HD3	5:B:659:HOH:O	2.20	0.42
1:A:422:GLN:HG2	1:A:423:LEU:HG	2.02	0.42
1:B:398:GLN:NE2	1:B:398:GLN:HA	2.35	0.42
1:B:107:LEU:HD23	1:B:107:LEU:HA	1.85	0.41
1:B:114:ARG:NH1	4:B:502:CHD:O26	2.53	0.41
4:B:502:CHD:H222	4:B:502:CHD:H162	1.91	0.41
1:A:352:TYR:HA	1:A:356:LEU:HB3	2.02	0.41
1:A:212:GLN:C	1:A:214:GLY:N	2.78	0.41
1:A:373:GLY:O	1:A:374:ASN:C	2.63	0.41
1:A:276:TYR:O	1:A:277:PRO:C	2.62	0.41
1:A:158:LYS:HD3	1:A:160:TYR:OH	2.21	0.41
1:A:305:VAL:HG23	4:A:504:CHD:H7	2.02	0.41
4:B:502:CHD:H183	4:B:502:CHD:H20	1.83	0.41
1:B:158:LYS:HD3	1:B:160:TYR:CZ	2.55	0.41
1:A:101:LEU:HB2	1:A:104:GLN:CG	2.51	0.40
1:A:304:LYS:NZ	1:A:308:MET:O	2.54	0.40
1:B:160:TYR:OH	1:B:182:LEU:HD11	2.22	0.40
1:B:207:TYR:CZ	1:B:413:GLU:HB3	2.57	0.40
1:B:335:ILE:HG13	1:B:335:ILE:O	2.21	0.40
1:A:160:TYR:OH	1:A:182:LEU:HD11	2.21	0.40
1:B:209:TYR:O	1:B:213:VAL:HG22	2.21	0.40
4:B:502:CHD:H183	4:B:504:CHD:H151	2.04	0.40
4:B:504:CHD:H21	4:B:504:CHD:H9	1.81	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	357/359 (99%)	331 (93%)	21 (6%)	5 (1%)	9	17
1	B	358/359 (100%)	328 (92%)	26 (7%)	4 (1%)	11	22
All	All	715/718 (100%)	659 (92%)	47 (7%)	9 (1%)	9	18

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	362	VAL
1	A	155	ALA
1	A	359	GLU
1	B	155	ALA
1	A	213	VAL
1	A	372	ASN
1	B	227	TRP
1	B	364	ASN
1	A	362	VAL

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	324/324 (100%)	298 (92%)	26 (8%)	11	24
1	B	325/324 (100%)	305 (94%)	20 (6%)	16	34
All	All	649/648 (100%)	603 (93%)	46 (7%)	13	29

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	96	ARG
1	A	98	LEU
1	A	100	THR
1	A	101	LEU

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Mol	Chain	Res	Type
1	A	103	ILE
1	A	124	ARG
1	A	132	ILE
1	A	153	ASN
1	A	160	TYR
1	A	183	GLU
1	A	211	ASN
1	A	212	GLN
1	A	213	VAL
1	A	220	LYS
1	A	292	GLU
1	A	298	ARG
1	A	341	HIS
1	A	342	ILE
1	A	358	LYS
1	A	367	ARG
1	A	374	ASN
1	A	378	SER
1	A	379	LYS
1	A	393	GLU
1	A	403	CYS
1	A	412	ARG
1	B	96	ARG
1	B	101	LEU
1	B	103	ILE
1	B	114	ARG
1	B	118	LYS
1	B	132	ILE
1	B	145	LYS
1	B	150	LEU
1	B	171	GLU
1	B	175	GLU
1	B	180	ASP
1	B	202	SER
1	B	246	ASP
1	B	286	LYS
1	B	342	ILE
1	B	358	LYS
1	B	392	ASN
1	B	393	GLU
1	B	410	VAL
1	B	423	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	122	GLN
1	A	211	ASN
1	A	235	GLN
1	A	240	HIS
1	A	302	GLN
1	A	329	ASN
1	A	364	ASN
1	A	374	ASN
1	A	386	HIS
1	A	388	HIS
1	A	398	GLN
1	A	421	GLN
1	B	212	GLN
1	B	235	GLN
1	B	240	HIS
1	B	263	HIS
1	B	302	GLN
1	B	374	ASN
1	B	386	HIS
1	B	388	HIS
1	B	392	ASN
1	B	398	GLN
1	B	421	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

8 ligands are modelled in this entry.

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
4	CHD	B	504	-	32,32,32	0.65	0	51,51,51	1.58	8 (15%)
3	GOL	A	502	-	5,5,5	0.34	0	5,5,5	0.34	0
4	CHD	A	503	-	32,32,32	0.78	1 (3%)	51,51,51	1.30	6 (11%)
4	CHD	B	502	-	32,32,32	0.76	1 (3%)	51,51,51	2.18	18 (35%)
4	CHD	B	503	-	32,32,32	0.82	0	51,51,51	2.06	17 (33%)
2	FES	B	501	1	0,4,4	-	-	-	-	-
4	CHD	A	504	-	32,32,32	0.84	0	51,51,51	1.66	13 (25%)
2	FES	A	501	1	0,4,4	-	-	-	-	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	CHD	B	504	-	-	7/9/74/74	0/4/4/4
3	GOL	A	502	-	-	1/4/4/4	-
4	CHD	B	503	-	-	7/9/74/74	0/4/4/4
4	CHD	B	502	-	-	5/9/74/74	0/4/4/4
4	CHD	A	503	-	-	5/9/74/74	0/4/4/4
2	FES	B	501	1	-	-	0/1/1/1
4	CHD	A	504	-	-	4/9/74/74	0/4/4/4
2	FES	A	501	1	-	-	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	502	CHD	C13-C14	-2.21	1.51	1.55
4	A	503	CHD	C4-C3	2.20	1.55	1.52

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	502	CHD	C6-C5-C10	5.36	118.37	112.66
4	B	503	CHD	C17-C13-C14	-4.80	95.30	100.11
4	B	502	CHD	C17-C13-C14	4.62	104.75	100.11
4	B	503	CHD	C13-C17-C20	-4.54	113.98	119.48
4	B	502	CHD	C19-C10-C5	-4.50	102.92	110.44
4	A	504	CHD	C16-C17-C20	4.38	118.80	112.18
4	B	503	CHD	C14-C13-C12	4.37	111.41	107.42
4	B	502	CHD	C9-C8-C7	-4.25	106.50	111.86
4	B	504	CHD	C6-C5-C4	-4.15	106.48	111.23
4	B	503	CHD	C16-C17-C20	4.14	118.45	112.18
4	B	503	CHD	C16-C17-C13	4.10	107.52	103.54
4	B	502	CHD	C9-C10-C5	3.83	113.83	108.51
4	A	504	CHD	C17-C13-C12	-3.77	114.27	117.67
4	B	502	CHD	C13-C17-C20	-3.67	115.04	119.48
4	B	502	CHD	C1-C2-C3	3.64	115.31	110.48
4	B	502	CHD	C14-C13-C12	-3.40	104.30	107.42
4	A	504	CHD	C13-C17-C20	-3.38	115.39	119.48
4	B	504	CHD	C1-C10-C9	-3.25	106.29	111.34
4	B	502	CHD	C11-C12-C13	3.23	114.55	111.26
4	B	504	CHD	C10-C9-C8	3.22	115.43	111.84
4	B	504	CHD	C9-C10-C5	3.10	112.82	108.51
4	A	503	CHD	C6-C7-C8	3.09	114.87	111.50
4	A	504	CHD	C15-C14-C8	3.07	122.57	118.36
4	A	504	CHD	C9-C8-C7	-3.06	108.00	111.86
4	B	503	CHD	C23-C22-C20	-3.04	108.78	114.46
4	B	504	CHD	C4-C5-C10	3.00	115.85	112.66
4	B	502	CHD	C9-C11-C12	2.96	118.16	114.29
4	B	502	CHD	C1-C10-C5	2.88	111.88	107.75
4	A	503	CHD	C2-C1-C10	2.79	117.45	112.74
4	A	504	CHD	C6-C7-C8	2.72	114.47	111.50
4	B	503	CHD	C14-C8-C7	-2.64	108.35	111.85
4	B	503	CHD	O12-C12-C11	2.63	114.46	109.12
4	B	502	CHD	C18-C13-C12	2.60	111.66	109.06
4	B	504	CHD	C14-C8-C9	-2.60	106.11	109.75
4	B	503	CHD	C15-C14-C13	2.58	106.05	103.54
4	A	504	CHD	C22-C20-C17	2.57	115.66	110.33
4	B	503	CHD	C6-C5-C4	-2.55	108.31	111.23
4	B	503	CHD	C21-C20-C17	-2.55	109.06	112.88
4	B	504	CHD	C17-C13-C14	2.54	102.66	100.11
4	B	504	CHD	C5-C4-C3	2.54	116.54	112.71
4	B	503	CHD	C5-C6-C7	2.48	117.35	114.40
4	A	504	CHD	C18-C13-C12	2.46	111.52	109.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	502	CHD	C4-C5-C10	2.37	115.18	112.66
4	A	504	CHD	C21-C20-C17	-2.31	109.42	112.88
4	B	502	CHD	C22-C20-C17	2.25	115.00	110.33
4	A	503	CHD	C19-C10-C9	-2.24	108.17	111.18
4	B	502	CHD	C13-C14-C8	-2.20	111.93	114.72
4	B	503	CHD	C11-C12-C13	-2.19	109.03	111.26
4	A	503	CHD	C1-C10-C9	2.16	114.70	111.34
4	A	504	CHD	C5-C6-C7	2.16	116.96	114.40
4	B	503	CHD	C11-C9-C8	2.15	114.07	110.89
4	B	503	CHD	C6-C5-C10	2.13	114.93	112.66
4	A	504	CHD	C4-C5-C10	2.10	114.89	112.66
4	B	502	CHD	C16-C17-C13	2.09	105.57	103.54
4	A	504	CHD	C19-C10-C1	-2.09	104.99	108.31
4	A	503	CHD	C15-C14-C13	2.08	105.56	103.54
4	B	502	CHD	C11-C9-C8	2.08	113.97	110.89
4	B	502	CHD	C11-C9-C10	2.07	115.81	113.70
4	B	503	CHD	C1-C2-C3	2.07	113.23	110.48
4	A	504	CHD	C1-C10-C5	2.07	110.72	107.75
4	B	503	CHD	C1-C10-C9	2.04	114.50	111.34
4	A	503	CHD	C19-C10-C1	-2.01	105.11	108.31

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	503	CHD	C13-C17-C20-C21
4	B	503	CHD	C13-C17-C20-C22
4	B	503	CHD	C16-C17-C20-C21
4	B	503	CHD	C16-C17-C20-C22
4	B	502	CHD	C16-C17-C20-C21
4	B	502	CHD	C13-C17-C20-C21
4	B	502	CHD	C13-C17-C20-C22
4	B	502	CHD	C16-C17-C20-C22
4	A	503	CHD	C17-C20-C22-C23
4	B	503	CHD	C17-C20-C22-C23
4	A	503	CHD	C21-C20-C22-C23
4	A	503	CHD	C20-C22-C23-C24
4	B	503	CHD	C20-C22-C23-C24
4	B	503	CHD	C21-C20-C22-C23
4	B	504	CHD	C20-C22-C23-C24
4	B	502	CHD	C20-C22-C23-C24
4	B	504	CHD	C16-C17-C20-C21

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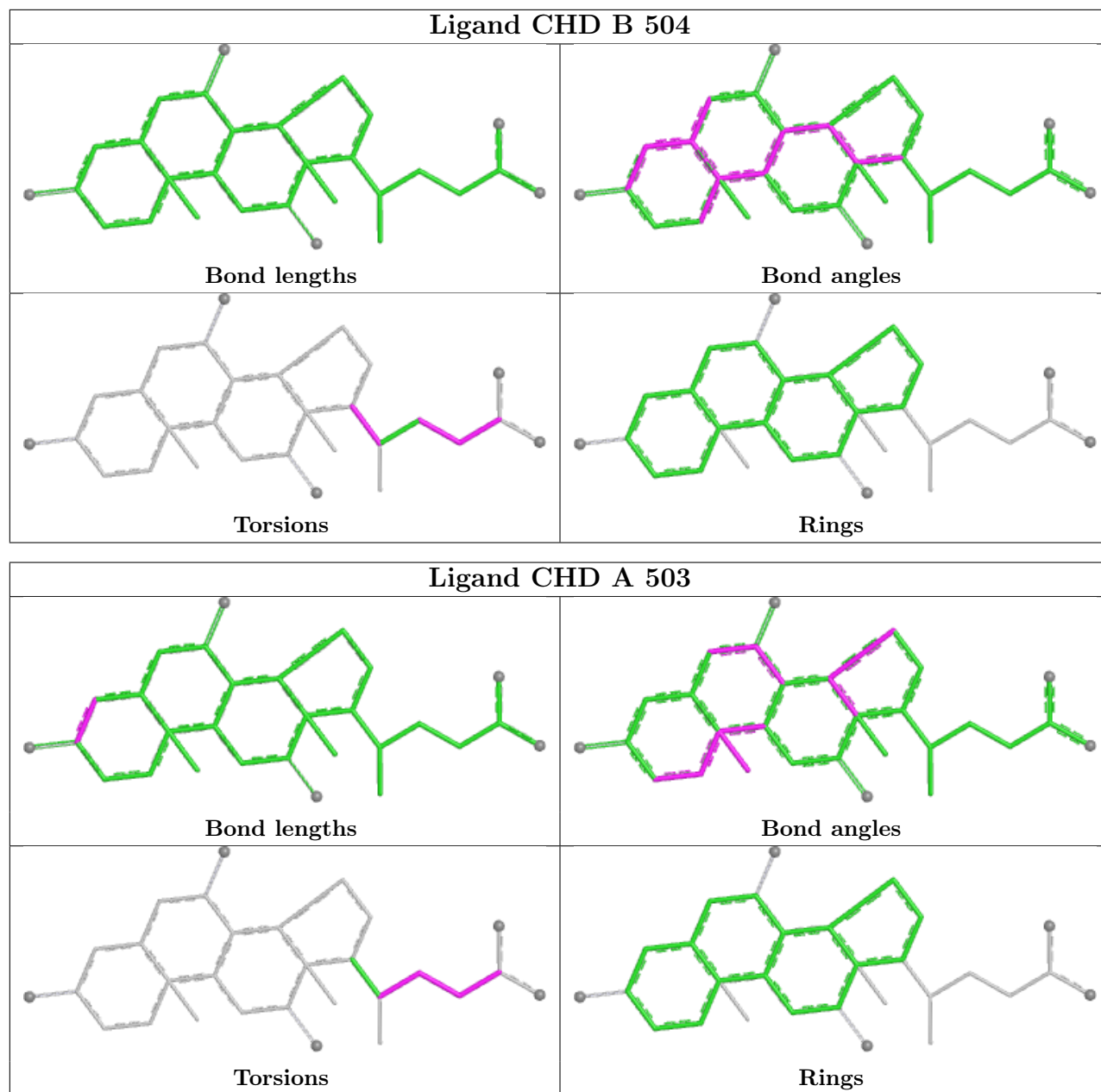
Mol	Chain	Res	Type	Atoms
4	B	504	CHD	C13-C17-C20-C22
4	B	504	CHD	C13-C17-C20-C21
4	B	504	CHD	C16-C17-C20-C22
4	A	504	CHD	C20-C22-C23-C24
3	A	502	GOL	O2-C2-C3-O3
4	A	504	CHD	C16-C17-C20-C22
4	A	503	CHD	C22-C23-C24-O25
4	A	503	CHD	C22-C23-C24-O26
4	B	504	CHD	C22-C23-C24-O26
4	B	504	CHD	C22-C23-C24-O25
4	A	504	CHD	C16-C17-C20-C21
4	A	504	CHD	C13-C17-C20-C21

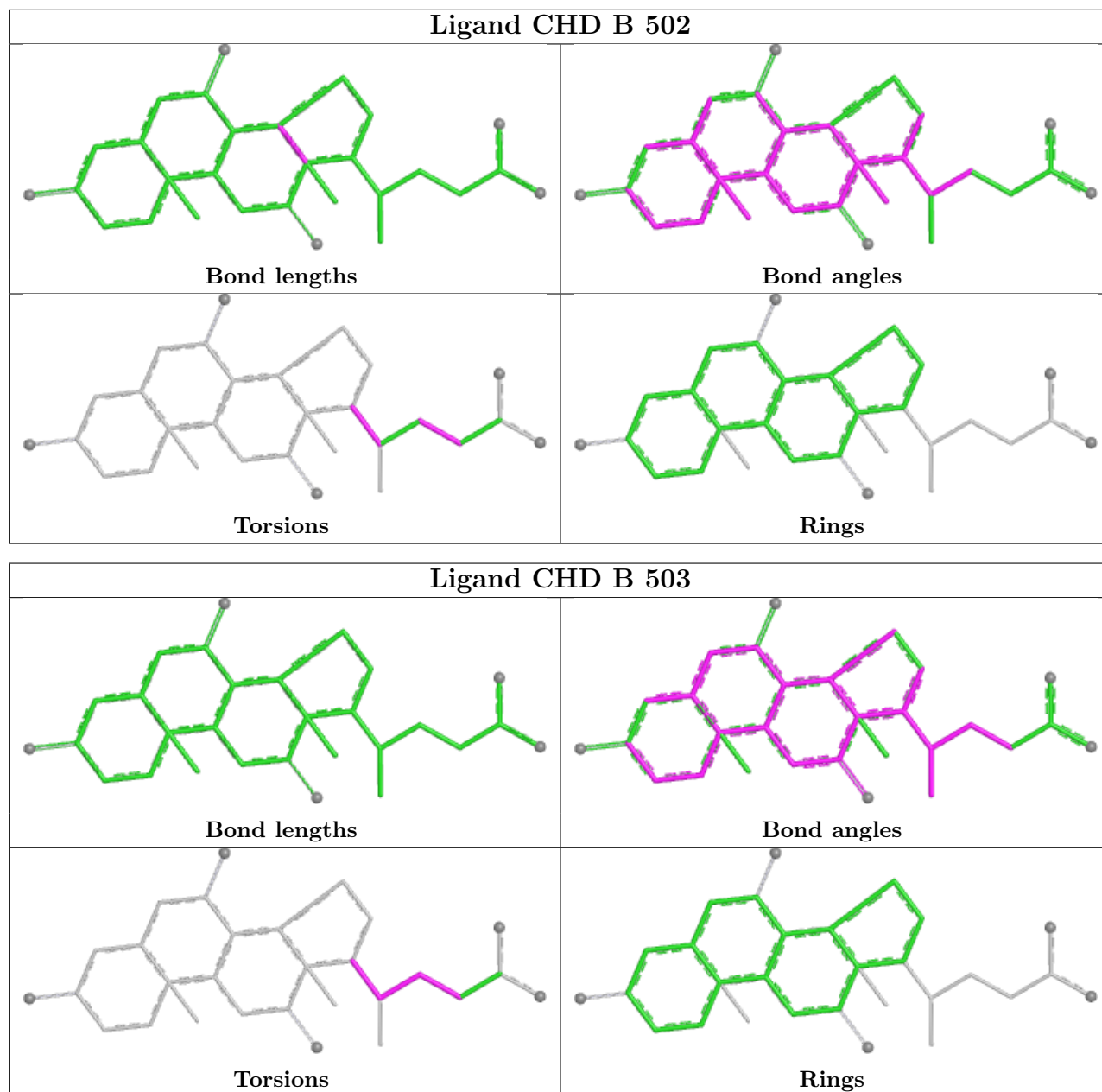
There are no ring outliers.

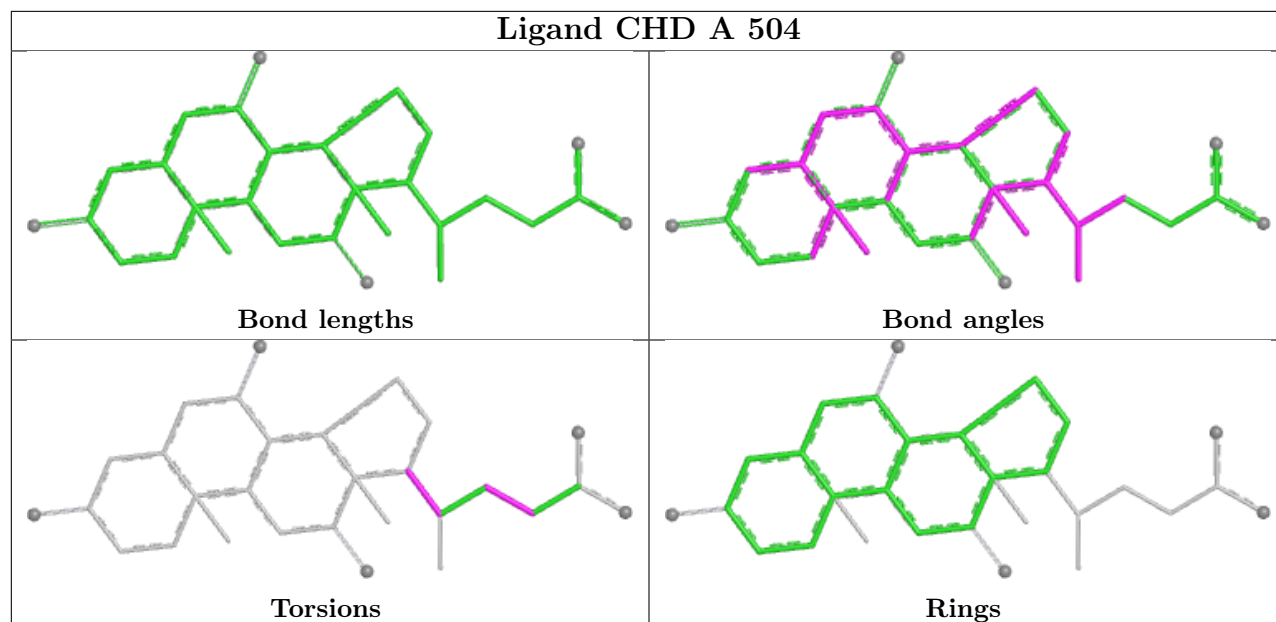
6 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	504	CHD	4	0
4	A	503	CHD	3	0
4	B	502	CHD	10	0
4	B	503	CHD	1	0
2	B	501	FES	1	0
4	A	504	CHD	3	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	359/359 (100%)	0.11	8 (2%) 62 58	18, 34, 55, 66	0
1	B	359/359 (100%)	0.13	11 (3%) 51 47	16, 33, 59, 69	1 (0%)
All	All	718/718 (100%)	0.12	19 (2%) 57 52	16, 33, 56, 69	1 (0%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	362	VAL	5.3
1	B	214	GLY	3.5
1	B	423	LEU	3.5
1	B	362	VAL	3.5
1	A	361	GLY	3.4
1	B	357	ALA	3.3
1	A	423	LEU	2.8
1	B	107	LEU	2.8
1	A	391	SER	2.7
1	A	76	MET	2.6
1	A	250	LEU	2.6
1	B	176	GLU	2.5
1	B	103	ILE	2.4
1	B	180	ASP	2.4
1	B	361	GLY	2.3
1	B	213	VAL	2.2
1	B	106	LYS	2.2
1	A	357	ALA	2.1
1	A	360	CYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

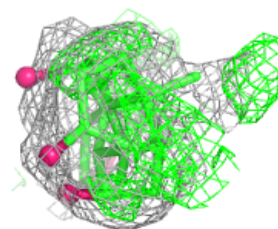
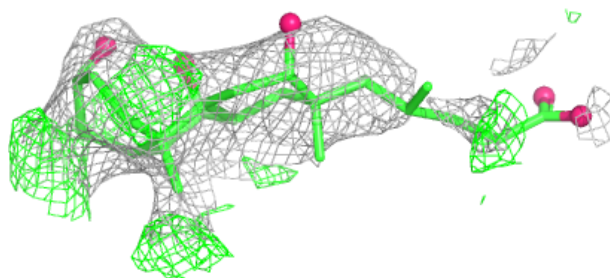
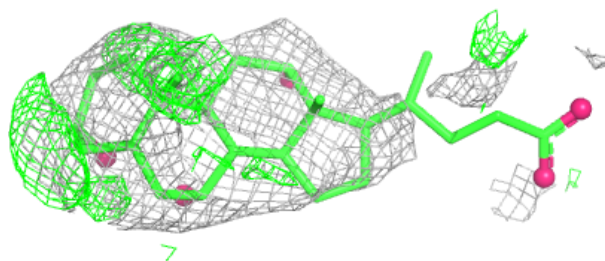
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	CHD	B	502	29/29	0.76	0.32	100,104,112,112	0
3	GOL	A	502	6/6	0.77	0.21	45,45,46,47	0
4	CHD	B	503	29/29	0.78	0.26	61,64,80,82	0
4	CHD	B	504	29/29	0.79	0.21	72,79,95,95	0
4	CHD	A	504	29/29	0.82	0.22	77,79,89,89	0
4	CHD	A	503	29/29	0.83	0.17	53,57,62,63	0
2	FES	A	501	4/4	0.99	0.05	36,39,41,43	0
2	FES	B	501	4/4	0.99	0.05	43,45,45,46	0

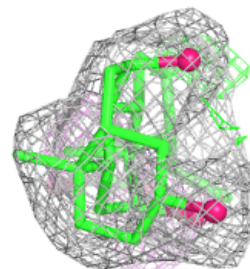
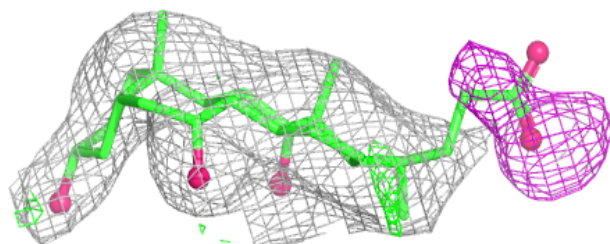
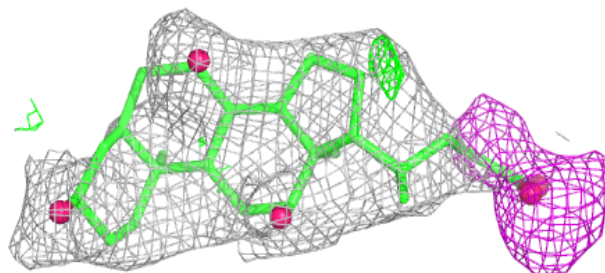
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around CHD B 502:

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

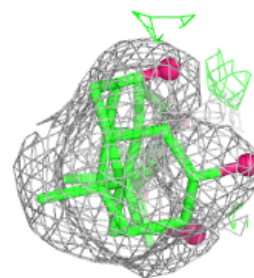
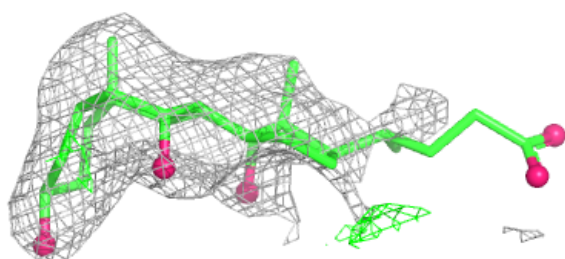
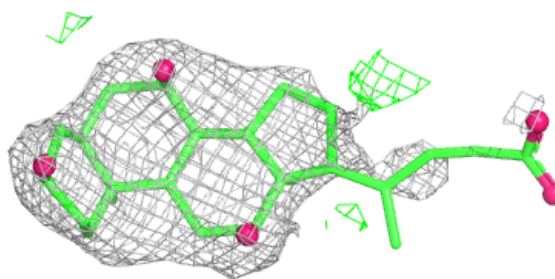
**Electron density around CHD B 503:**

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

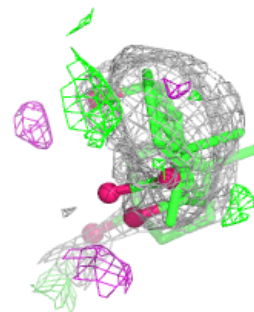
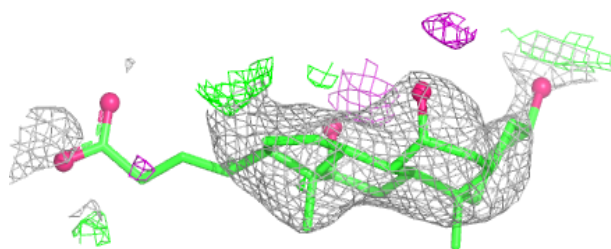
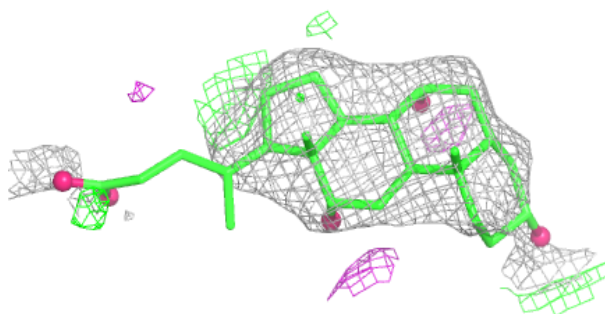


Electron density around CHD B 504:

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

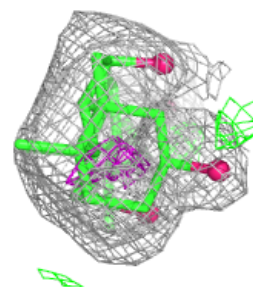
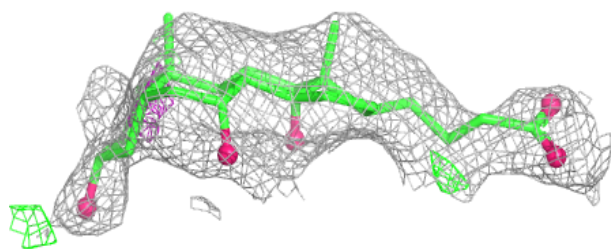
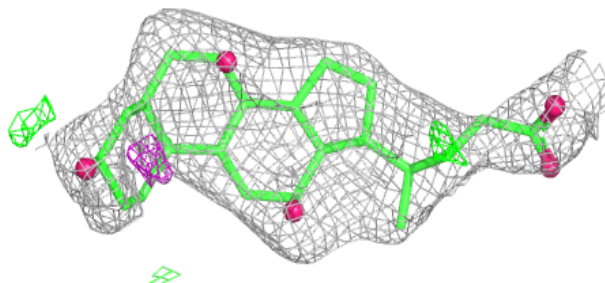
**Electron density around CHD A 504:**

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



Electron density around CHD A 503:

$2mF_o-DF_c$ (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.