



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:40 PM UTC

PDB ID : 6TXI / pdb_00006txi
Title : Crystal structure of thermotoga maritima E65A Ferritin
Authors : Wilk, P.; Grudnik, P.; Kumar, M.; Heddle, J.; Chakraborti, S.
Deposited on : 2020-01-14
Resolution : 1.76 Å(reported)

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

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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 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

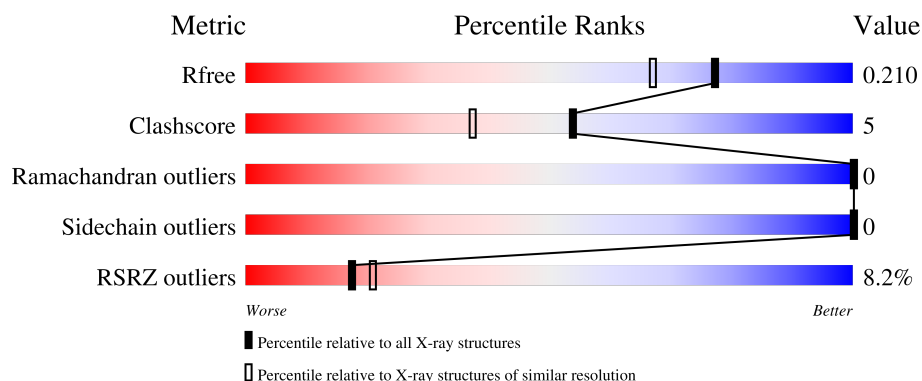
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.76 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.



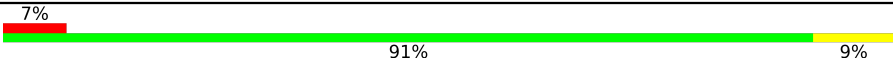
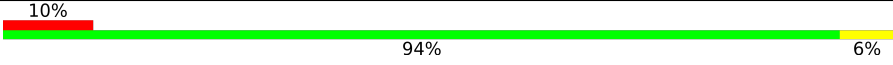
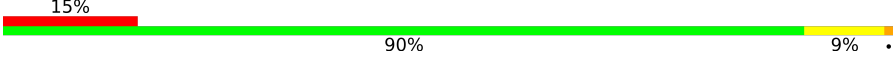
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (1.76-1.76)

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	164	5% 95% 5%
1	B	164	7% 93% 7%
1	C	164	9% 87% 10% ..
1	D	164	9% 91% 9% .
1	E	164	4% 92% 7% .

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Mol	Chain	Length	Quality of chain
1	F	164	
1	G	164	
1	H	164	

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
2	SO4	F	201	-	-	X	-
3	GOL	F	202	-	-	X	-
4	FE	D	204	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 23189 atoms, of which 10810 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 Ferritin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	164	Total	C	H	N	O	S	0	3	0
			2725	884	1345	226	264	6			
1	B	164	Total	C	H	N	O	S	0	0	0
			2681	872	1318	223	262	6			
1	C	164	Total	C	H	N	O	S	0	1	0
			2693	876	1324	223	264	6			
1	D	164	Total	C	H	N	O	S	0	2	0
			2711	880	1337	225	263	6			
1	E	164	Total	C	H	N	O	S	0	0	0
			2681	872	1318	223	262	6			
1	F	164	Total	C	H	N	O	S	0	0	0
			2681	872	1318	223	262	6			
1	G	164	Total	C	H	N	O	S	0	0	0
			2681	872	1318	223	262	6			
1	H	164	Total	C	H	N	O	S	0	1	0
			2700	877	1331	224	262	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	65	ALA	GLU	engineered mutation	UNP Q9X0L2
B	65	ALA	GLU	engineered mutation	UNP Q9X0L2
C	65	ALA	GLU	engineered mutation	UNP Q9X0L2
D	65	ALA	GLU	engineered mutation	UNP Q9X0L2
E	65	ALA	GLU	engineered mutation	UNP Q9X0L2
F	65	ALA	GLU	engineered mutation	UNP Q9X0L2
G	65	ALA	GLU	engineered mutation	UNP Q9X0L2
H	65	ALA	GLU	engineered mutation	UNP Q9X0L2

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			9	3	3	3		
3	B	1	Total	C	H	O	0	0
			9	3	3	3		
3	B	1	Total	C	H	O	0	0
			9	3	3	3		
3	C	1	Total	C	H	O	0	0
			9	3	3	3		
3	C	1	Total	C	H	O	0	0
			9	3	3	3		
3	C	1	Total	C	H	O	0	0
			9	3	3	3		
3	D	1	Total	C	H	O	0	0
			9	3	3	3		
3	E	1	Total	C	H	O	0	0
			9	3	3	3		
3	F	1	Total	C	H	O	0	0
			9	3	3	3		

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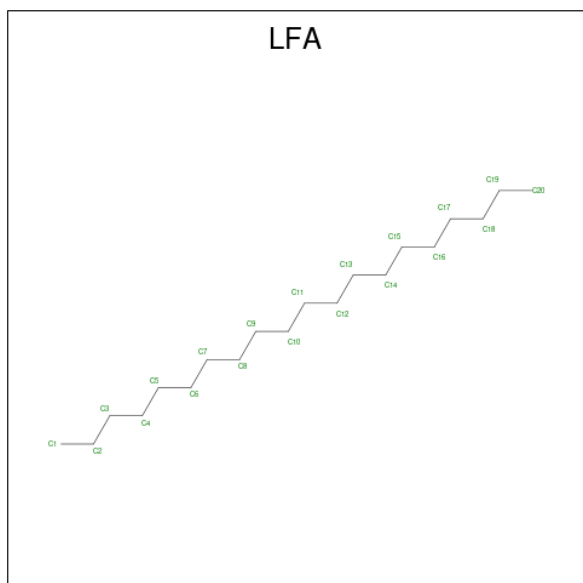
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	G	1	Total	C	H	O	0	0
			9	3	3	3		
3	G	1	Total	C	H	O	0	0
			9	3	3	3		

- Molecule 4 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Fe	0	0
			1	1		
4	B	1	Total	Fe	0	0
			1	1		
4	C	1	Total	Fe	0	0
			1	1		
4	D	1	Total	Fe	0	0
			1	1		
4	E	1	Total	Fe	0	0
			1	1		
4	F	1	Total	Fe	0	0
			1	1		
4	G	1	Total	Fe	0	0
			1	1		
4	H	1	Total	Fe	0	0
			1	1		

- Molecule 5 is EICOSANE (CCD ID: LFA) (formula: C₂₀H₄₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total C H 62 20 42	0	0
5	C	1	Total C H 62 20 42	0	0
5	F	1	Total C H 62 20 42	0	0
5	G	1	Total C H 62 20 42	0	0

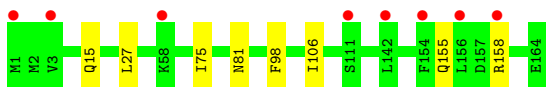
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	176	Total O 176 176	0	0
6	B	190	Total O 190 190	0	0
6	C	134	Total O 134 134	0	0
6	D	160	Total O 160 160	0	0
6	E	163	Total O 163 163	0	0
6	F	147	Total O 147 147	0	0
6	G	122	Total O 122 122	0	0
6	H	109	Total O 109 109	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.

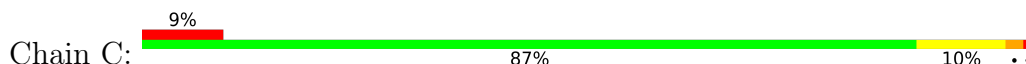
- Molecule 1: Ferritin



- Molecule 1: Ferritin



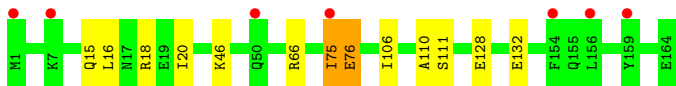
- Molecule 1: Ferritin



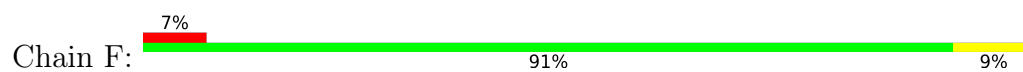
- Molecule 1: Ferritin



- Molecule 1: Ferritin



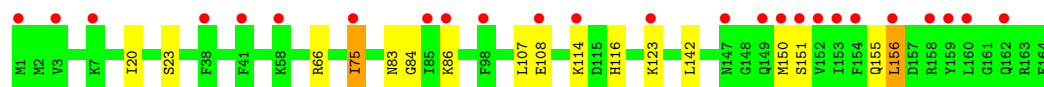
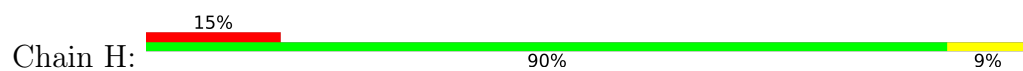
- Molecule 1: Ferritin



• Molecule 1: Ferritin



• Molecule 1: Ferritin



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	175.70Å 175.70Å 353.57Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.94 – 1.76 48.94 – 1.76	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.94-1.76) 99.9 (48.94-1.76)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.99 (at 1.76Å)	Xtriage
Refinement program	PHENIX 1.15.1	Depositor
R, R_{free}	0.180 , 0.204 0.185 , 0.210	Depositor DCC
R_{free} test set	2168 reflections (1.05%)	wwPDB-VP
Wilson B-factor (Å ²)	37.3	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 61.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	23189	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.61% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, LFA, FE, GOL

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	0.44	0/1418	0.56	0/1905
1	B	0.40	0/1392	0.57	1/1871 (0.1%)
1	C	0.78	6/1401 (0.4%)	0.92	9/1883 (0.5%)
1	D	0.42	0/1409	0.62	2/1893 (0.1%)
1	E	0.56	2/1392 (0.1%)	0.70	2/1871 (0.1%)
1	F	0.36	0/1392	0.53	0/1871
1	G	0.39	0/1392	0.62	2/1871 (0.1%)
1	H	0.70	5/1401 (0.4%)	0.87	8/1882 (0.4%)
All	All	0.53	13/11197 (0.1%)	0.69	24/15047 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	2
1	D	0	1
All	All	0	3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	123	LYS	CE-NZ	-10.22	1.18	1.49
1	C	135	GLN	CD-NE2	9.21	1.52	1.33
1	C	135	GLN	CB-CG	8.99	1.79	1.52
1	H	84	GLY	C-O	7.84	1.34	1.23
1	C	144	GLU	CD-OE2	7.26	1.39	1.25
1	C	135	GLN	CA-CB	6.52	1.63	1.53
1	C	135	GLN	CG-CD	6.19	1.67	1.52
1	E	110	ALA	C-O	-5.90	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	86	LYS	CE-NZ	5.89	1.67	1.49
1	H	123	LYS	CG-CD	5.67	1.69	1.52
1	C	85	ILE	CG1-CD1	5.28	1.72	1.51
1	E	111	SER	C-O	-5.16	1.18	1.24
1	H	114	LYS	CD-CE	5.15	1.67	1.52

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	135	GLN	CA-CB-CG	11.83	137.76	114.10
1	C	158	ARG	NE-CZ-NH2	11.59	129.63	119.20
1	C	158	ARG	CG-CD-NE	10.97	136.12	112.00
1	H	86	LYS	CD-CE-NZ	-9.87	80.32	111.90
1	H	86	LYS	CG-CD-CE	9.35	132.81	111.30
1	H	156	LEU	CA-CB-CG	-8.85	85.31	116.30
1	H	114	LYS	CB-CG-CD	-8.74	91.20	111.30
1	E	75	ILE	CA-CB-CG1	8.26	124.43	110.40
1	C	153	ILE	CA-CB-CG1	-7.58	97.52	110.40
1	C	86	LYS	CD-CE-NZ	-7.25	88.70	111.90
1	C	158	ARG	CD-NE-CZ	7.21	134.49	124.40
1	H	114	LYS	CD-CE-NZ	-6.99	89.54	111.90
1	G	158	ARG	CG-CD-NE	-6.86	96.90	112.00
1	C	158	ARG	NE-CZ-NH1	-6.86	114.64	121.50
1	H	75	ILE	CA-CB-CG1	6.82	122.00	110.40
1	D	149	GLN	CA-CB-CG	6.19	126.48	114.10
1	D	153	ILE	CA-CB-CG1	-6.04	100.13	110.40
1	H	114	LYS	N-CA-CB	-6.03	103.30	112.47
1	G	158	ARG	CB-CA-C	-6.00	101.45	110.88
1	C	85	ILE	CA-CB-CG1	5.78	120.22	110.40
1	H	114	LYS	CA-CB-CG	5.71	125.52	114.10
1	C	135	GLN	N-CA-CB	5.19	117.84	110.16
1	E	76	GLU	CG-CD-OE2	5.07	130.06	118.40
1	B	18	ARG	CG-CD-NE	-5.03	100.94	112.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	135	GLN	Sidechain
1	C	158	ARG	Sidechain
1	D	149	GLN	Sidechain

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	1380	1345	1345	9	0
1	B	1363	1318	1318	10	0
1	C	1369	1324	1324	20	0
1	D	1374	1337	1337	11	0
1	E	1363	1318	1318	14	0
1	F	1363	1318	1318	18	0
1	G	1363	1318	1318	11	0
1	H	1369	1331	1331	13	2
2	A	15	0	0	1	0
2	B	20	0	0	2	0
2	C	5	0	0	1	0
2	D	10	0	0	1	0
2	F	10	0	0	3	0
2	G	5	0	0	0	0
2	H	15	0	0	0	0
3	A	6	3	8	1	0
3	B	12	6	16	2	0
3	C	18	9	24	5	0
3	D	6	3	8	0	0
3	E	6	3	8	1	0
3	F	6	3	8	6	0
3	G	12	6	16	3	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0
5	B	20	42	42	2	0
5	C	20	42	42	2	0
5	F	20	42	42	2	0
5	G	20	42	42	3	0
6	A	176	0	0	2	0
6	B	190	0	0	4	0
6	C	134	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	D	160	0	0	3	0
6	E	163	0	0	1	0
6	F	147	0	0	4	0
6	G	122	0	0	1	0
6	H	109	0	0	2	0
All	All	12379	10810	10865	107	2

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 5.

All (107) 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:135:GLN:CB	1:C:135:GLN:CG	1.79	1.58
1:C:18:ARG:HH11	3:C:202:GOL:H31	1.13	1.14
1:C:18:ARG:NH1	3:C:202:GOL:H31	1.80	0.95
2:F:201:SO4:O4	6:F:301:HOH:O	1.92	0.85
2:A:204:SO4:O3	6:A:301:HOH:O	1.95	0.85
1:A:155[B]:GLN:CD	1:A:158:ARG:HH12	1.88	0.80
2:C:206:SO4:O3	6:C:301:HOH:O	1.98	0.80
2:B:201:SO4:O1	6:B:301:HOH:O	2.02	0.77
2:F:204:SO4:O3	6:F:302:HOH:O	2.02	0.76
1:A:98:PHE:HZ	3:A:202:GOL:H12	1.52	0.73
1:F:98:PHE:CZ	3:F:202:GOL:H11	2.24	0.72
1:F:98:PHE:HZ	3:F:202:GOL:H11	1.54	0.72
1:D:50:GLN:NE2	6:D:302:HOH:O	2.25	0.70
1:H:142:LEU:HD13	1:H:156:LEU:HD21	1.73	0.69
2:F:201:SO4:O2	6:F:303:HOH:O	2.11	0.66
1:B:104:TYR:OH	6:B:302:HOH:O	2.12	0.65
2:B:207:SO4:O2	6:B:303:HOH:O	2.14	0.62
1:E:46:LYS:HB3	5:F:203:LFA:H21	1.82	0.62
1:G:18:ARG:HE	3:G:203:GOL:H2	1.66	0.61
1:A:155[B]:GLN:OE1	1:A:158:ARG:NH1	2.34	0.61
1:E:18:ARG:NH2	3:E:202:GOL:O2	2.21	0.61
3:C:204:GOL:H12	6:C:340:HOH:O	2.01	0.60
1:G:86:LYS:H	3:G:202:GOL:H2	1.66	0.60
1:E:75:ILE:HB	1:F:75:ILE:HB	1.83	0.60
1:C:41:PHE:CD2	1:C:85:ILE:HD12	2.37	0.59
1:C:33:PHE:CE1	1:C:85:ILE:HD13	2.39	0.58
1:A:155[B]:GLN:CD	1:A:158:ARG:NH1	2.61	0.57
1:G:75:ILE:HB	1:H:75:ILE:HB	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:83:ASN:ND2	6:H:302:HOH:O	2.25	0.57
1:C:86:LYS:HE3	1:C:144:GLU:OE1	2.05	0.57
1:F:18:ARG:HG2	3:F:202:GOL:H12	1.85	0.57
1:D:146:ALA:O	1:D:149:GLN:HB2	2.05	0.56
1:C:41:PHE:CD1	1:C:160:LEU:HD12	2.41	0.56
1:B:46:LYS:HB3	5:B:206:LFA:H32	1.87	0.56
1:F:41:PHE:CD1	1:F:160:LEU:HD12	2.40	0.56
1:E:75:ILE:HG13	1:F:75:ILE:HB	1.89	0.54
1:C:3:VAL:O	3:C:204:GOL:H11	2.07	0.54
1:A:27:LEU:HD22	5:B:206:LFA:H171	1.89	0.53
1:F:38:PHE:HB3	1:F:41:PHE:HD2	1.73	0.53
1:E:20:ILE:HG21	1:E:75:ILE:HD11	1.91	0.53
1:D:41:PHE:CD1	1:D:160:LEU:HD12	2.44	0.52
1:E:20:ILE:HG22	1:E:75:ILE:CD1	2.40	0.52
1:C:57:MET:HE1	1:C:60:TYR:CD2	2.44	0.52
1:C:86:LYS:HE2	1:C:90:GLU:OE2	2.10	0.52
1:C:155:GLN:NE2	6:C:302:HOH:O	2.44	0.51
1:G:41:PHE:CD1	1:G:160:LEU:HD12	2.45	0.51
1:A:155[B]:GLN:NE2	1:A:158:ARG:HH12	2.09	0.51
1:C:18:ARG:HD3	3:C:202:GOL:O2	2.11	0.51
1:G:50:GLN:HG2	5:G:204:LFA:H181	1.94	0.50
1:F:41:PHE:CE1	1:F:160:LEU:HD12	2.47	0.50
1:E:20:ILE:CG2	1:E:75:ILE:CD1	2.90	0.50
1:H:20:ILE:HG21	1:H:75:ILE:HD11	1.93	0.50
1:D:41:PHE:HD1	1:D:160:LEU:HD12	1.77	0.49
5:C:205:LFA:H22	1:D:50:GLN:OE1	2.11	0.49
1:H:155:GLN:C	1:H:156:LEU:HD12	2.37	0.49
1:E:75:ILE:CG1	1:F:75:ILE:HB	2.42	0.49
1:C:135:GLN:CB	1:C:135:GLN:CD	2.79	0.48
1:C:41:PHE:HD1	1:C:160:LEU:HD12	1.78	0.48
1:G:75:ILE:HB	1:H:75:ILE:HG13	1.96	0.48
1:E:15:GLN:HB2	1:E:106:ILE:HD11	1.95	0.48
1:G:41:PHE:CE1	1:G:160:LEU:HD12	2.49	0.48
1:D:86:LYS:HB2	2:D:201:SO4:O4	2.14	0.47
1:H:142:LEU:HD13	1:H:156:LEU:CD2	2.42	0.47
1:D:58:LYS:NZ	6:D:305:HOH:O	2.32	0.47
1:A:15:GLN:HB2	1:A:106:ILE:HD11	1.96	0.47
1:B:10:LYS:HD2	6:B:309:HOH:O	2.15	0.47
1:B:15:GLN:HB2	1:B:106:ILE:HD11	1.97	0.47
1:H:20:ILE:HG22	1:H:75:ILE:CD1	2.45	0.46
1:D:103:ILE:HD13	1:D:125:PHE:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:15:GLN:HB2	1:G:106:ILE:HD11	1.96	0.46
1:G:50:GLN:HE21	5:G:204:LFA:H203	1.80	0.46
1:A:81[B]:ASN:ND2	6:A:309:HOH:O	2.50	0.45
5:G:204:LFA:H72	1:H:23:SER:OG	2.16	0.45
1:F:15:GLN:HB2	1:F:106:ILE:HD11	1.98	0.45
1:F:86:LYS:O	1:F:90:GLU:HG2	2.17	0.45
1:F:18:ARG:NE	3:F:202:GOL:H12	2.32	0.44
1:C:15:GLN:HB2	1:C:106:ILE:HD11	1.99	0.44
1:C:33:PHE:HE1	1:C:85:ILE:HD13	1.82	0.44
1:F:132:GLU:O	1:F:135:GLN:HG2	2.18	0.44
1:H:20:ILE:CG2	1:H:75:ILE:CD1	2.95	0.44
1:B:98:PHE:HZ	3:B:205:GOL:H32	1.83	0.43
1:B:98:PHE:CZ	3:B:205:GOL:H32	2.53	0.43
1:G:86:LYS:HB2	3:G:202:GOL:H12	2.00	0.43
1:F:7:LYS:CE	6:F:318:HOH:O	2.65	0.43
1:H:66:ARG:HA	1:H:66:ARG:HD2	1.87	0.43
1:B:160:LEU:N	1:B:160:LEU:HD23	2.33	0.43
1:F:57:MET:HG3	5:F:203:LFA:H32	1.99	0.43
1:E:128:GLU:O	1:E:132:GLU:HG2	2.19	0.43
1:A:75:ILE:HB	1:B:75:ILE:HB	2.01	0.43
1:B:128:GLU:O	1:B:132:GLU:HG2	2.19	0.43
1:C:38:PHE:CD2	1:C:85:ILE:HD11	2.54	0.42
1:G:90:GLU:OE1	6:G:301:HOH:O	2.22	0.42
1:E:16:LEU:C	1:E:16:LEU:HD23	2.45	0.42
1:F:18:ARG:CD	3:F:202:GOL:H12	2.50	0.42
1:F:18:ARG:CG	3:F:202:GOL:H12	2.49	0.42
1:H:150:MET:O	1:H:151:SER:C	2.62	0.42
5:C:205:LFA:H201	6:D:460:HOH:O	2.20	0.41
1:H:107:LEU:HD11	6:H:301:HOH:O	2.20	0.41
1:E:75:ILE:CB	1:F:75:ILE:HB	2.49	0.41
1:C:119:VAL:HG12	1:C:123:LYS:HE3	2.02	0.41
1:D:15:GLN:HB2	1:D:106:ILE:HD11	2.02	0.41
1:B:132:GLU:O	1:B:135:GLN:HG2	2.21	0.41
1:E:76:GLU:HG3	6:E:408:HOH:O	2.21	0.41
1:C:35:ALA:HB2	1:D:69:ARG:HG3	2.03	0.40
1:C:158:ARG:O	1:C:162:GLN:HG3	2.20	0.40
1:D:107:LEU:HD11	1:D:123[A]:LYS:HE2	2.03	0.40
1:E:66:ARG:HA	1:E:66:ARG:HD2	1.90	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:108:GLU:OE2	1:H:116:HIS:HE2[3_565]	1.24	0.36
1:H:108:GLU:OE2	1:H:116:HIS:NE2[3_565]	2.00	0.20

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	165/164 (101%)	165 (100%)	0	0	100	100
1	B	162/164 (99%)	162 (100%)	0	0	100	100
1	C	163/164 (99%)	162 (99%)	1 (1%)	0	100	100
1	D	164/164 (100%)	163 (99%)	1 (1%)	0	100	100
1	E	162/164 (99%)	162 (100%)	0	0	100	100
1	F	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	G	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	H	163/164 (99%)	162 (99%)	1 (1%)	0	100	100
All	All	1303/1312 (99%)	1298 (100%)	5 (0%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	146/143 (102%)	146 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	143/143 (100%)	143 (100%)	0	100	100
1	C	144/143 (101%)	144 (100%)	0	100	100
1	D	145/143 (101%)	145 (100%)	0	100	100
1	E	143/143 (100%)	143 (100%)	0	100	100
1	F	143/143 (100%)	143 (100%)	0	100	100
1	G	143/143 (100%)	143 (100%)	0	100	100
1	H	144/143 (101%)	144 (100%)	0	100	100
All	All	1151/1144 (101%)	1151 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	129	GLN
1	C	129	GLN
1	D	101	GLN
1	D	116	HIS
1	D	129	GLN
1	D	149	GLN
1	E	129	GLN
1	E	155	GLN
1	G	50	GLN
1	G	129	GLN
1	H	81	ASN
1	H	83	ASN
1	H	129	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 39 ligands modelled in this entry, 8 are monoatomic - leaving 31 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	$\# Z > 2$
2	SO4	B	202	-	4,4,4	0.72	0	6,6,6	0.14	0
3	GOL	F	202	-	5,5,5	0.85	0	5,5,5	1.02	0
3	GOL	A	202	-	5,5,5	0.23	0	5,5,5	0.29	0
5	LFA	B	206	-	19,19,19	0.56	0	18,18,18	0.26	0
2	SO4	A	204	-	4,4,4	0.21	0	6,6,6	0.12	0
5	LFA	C	205	-	19,19,19	0.39	0	18,18,18	0.30	0
5	LFA	G	204	-	19,19,19	0.58	0	18,18,18	0.26	0
3	GOL	B	205	-	5,5,5	0.86	0	5,5,5	1.20	1 (20%)
2	SO4	A	201	-	4,4,4	0.54	0	6,6,6	0.16	0
3	GOL	G	202	-	5,5,5	0.78	0	5,5,5	0.94	0
3	GOL	D	202	-	5,5,5	0.16	0	5,5,5	0.23	0
2	SO4	B	207	-	4,4,4	0.24	0	6,6,6	0.08	0
2	SO4	D	201	-	4,4,4	0.74	0	6,6,6	0.26	0
2	SO4	D	203	-	4,4,4	0.24	0	6,6,6	0.15	0
2	SO4	A	203	-	4,4,4	0.29	0	6,6,6	0.12	0
2	SO4	G	205	-	4,4,4	0.32	0	6,6,6	0.06	0
3	GOL	C	203	-	5,5,5	0.18	0	5,5,5	0.32	0
2	SO4	B	204	-	4,4,4	1.26	0	6,6,6	0.10	0
2	SO4	F	204	-	4,4,4	0.24	0	6,6,6	0.07	0
3	GOL	C	202	-	5,5,5	0.21	0	5,5,5	0.58	0
3	GOL	C	204	-	5,5,5	0.18	0	5,5,5	0.43	0
2	SO4	F	201	-	4,4,4	0.73	0	6,6,6	0.12	0
2	SO4	H	203	-	4,4,4	0.26	0	6,6,6	0.12	0
3	GOL	E	202	-	5,5,5	0.23	0	5,5,5	0.43	0
2	SO4	H	201	-	4,4,4	0.52	0	6,6,6	0.18	0
2	SO4	B	201	-	4,4,4	0.65	0	6,6,6	0.12	0
3	GOL	B	203	-	5,5,5	0.26	0	5,5,5	0.21	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	C	206	-	4,4,4	0.23	0	6,6,6	0.08	0
5	LFA	F	203	-	19,19,19	0.52	0	18,18,18	0.13	0
3	GOL	G	203	-	5,5,5	0.18	0	5,5,5	0.33	0
2	SO4	H	202	-	4,4,4	0.27	0	6,6,6	0.11	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
3	GOL	B	203	-	-	0/4/4/4	-
3	GOL	F	202	-	-	4/4/4/4	-
5	LFA	F	203	-	-	1/17/17/17	-
3	GOL	C	203	-	-	0/4/4/4	-
3	GOL	G	203	-	-	2/4/4/4	-
3	GOL	B	205	-	-	3/4/4/4	-
3	GOL	C	202	-	-	2/4/4/4	-
3	GOL	C	204	-	-	0/4/4/4	-
3	GOL	A	202	-	-	2/4/4/4	-
3	GOL	G	202	-	-	2/4/4/4	-
3	GOL	D	202	-	-	0/4/4/4	-
5	LFA	B	206	-	-	2/17/17/17	-
3	GOL	E	202	-	-	3/4/4/4	-
5	LFA	C	205	-	-	5/17/17/17	-
5	LFA	G	204	-	-	6/17/17/17	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	205	GOL	C3-C2-C1	-2.19	103.75	111.80

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	F	202	GOL	O1-C1-C2-O2
3	F	202	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
3	G	203	GOL	O1-C1-C2-C3
3	A	202	GOL	C1-C2-C3-O3
3	B	205	GOL	O1-C1-C2-C3
3	C	202	GOL	C1-C2-C3-O3
3	E	202	GOL	O1-C1-C2-C3
3	F	202	GOL	C1-C2-C3-O3
3	G	202	GOL	C1-C2-C3-O3
3	C	202	GOL	O2-C2-C3-O3
3	G	202	GOL	O2-C2-C3-O3
3	G	203	GOL	O1-C1-C2-O2
5	G	204	LFA	C2-C3-C4-C5
3	B	205	GOL	O1-C1-C2-O2
5	C	205	LFA	C10-C11-C12-C13
5	G	204	LFA	C11-C10-C9-C8
5	C	205	LFA	C12-C13-C14-C15
5	C	205	LFA	C11-C12-C13-C14
5	C	205	LFA	C9-C10-C11-C12
3	B	205	GOL	O2-C2-C3-O3
3	F	202	GOL	O2-C2-C3-O3
5	G	204	LFA	C16-C17-C18-C19
5	G	204	LFA	C6-C7-C8-C9
3	E	202	GOL	O1-C1-C2-O2
3	E	202	GOL	O2-C2-C3-O3
5	B	206	LFA	C11-C10-C9-C8
5	F	203	LFA	C7-C8-C9-C10
5	B	206	LFA	C1-C2-C3-C4
3	A	202	GOL	O2-C2-C3-O3
5	G	204	LFA	C7-C8-C9-C10
5	C	205	LFA	C7-C8-C9-C10
5	G	204	LFA	C11-C12-C13-C14

There are no ring outliers.

19 monomers are involved in 35 short contacts:

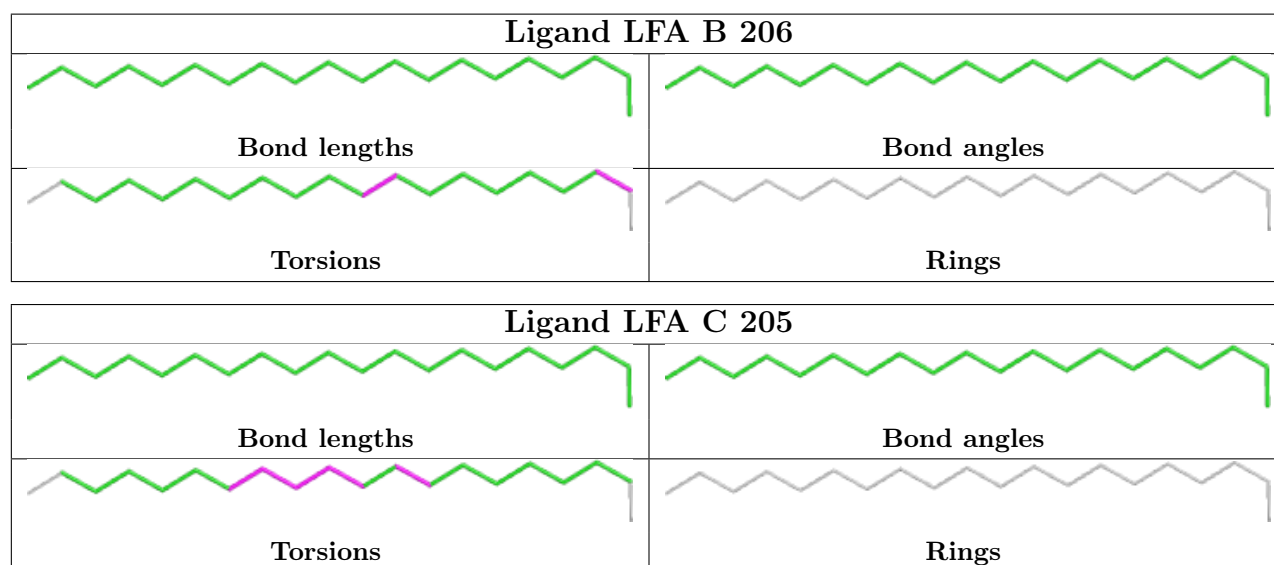
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	202	GOL	6	0
3	A	202	GOL	1	0
5	B	206	LFA	2	0
2	A	204	SO4	1	0
5	C	205	LFA	2	0
5	G	204	LFA	3	0
3	B	205	GOL	2	0

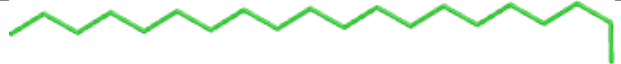
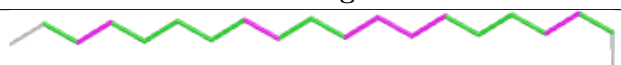
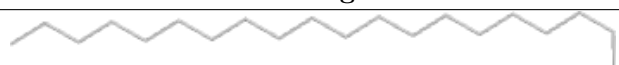
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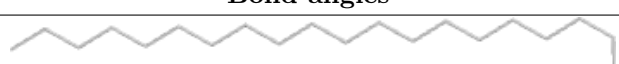
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	202	GOL	2	0
2	B	207	SO4	1	0
2	D	201	SO4	1	0
2	F	204	SO4	1	0
3	C	202	GOL	3	0
3	C	204	GOL	2	0
2	F	201	SO4	2	0
3	E	202	GOL	1	0
2	B	201	SO4	1	0
2	C	206	SO4	1	0
5	F	203	LFA	2	0
3	G	203	GOL	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



Ligand LFA G 204	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand LFA F 203	
 Bond lengths	 Bond angles
 Torsions	 Rings

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	164/164 (100%)	0.23	8 (4%)	35 40	24, 43, 75, 135	3 (1%)
1	B	164/164 (100%)	0.23	12 (7%)	21 24	32, 41, 78, 119	0
1	C	164/164 (100%)	0.46	14 (8%)	16 20	37, 50, 92, 113	1 (0%)
1	D	164/164 (100%)	0.37	14 (8%)	16 20	23, 46, 80, 143	2 (1%)
1	E	164/164 (100%)	0.18	7 (4%)	40 46	33, 44, 84, 134	0
1	F	164/164 (100%)	0.27	11 (6%)	24 28	35, 48, 81, 117	0
1	G	164/164 (100%)	0.59	17 (10%)	11 13	42, 55, 90, 125	0
1	H	164/164 (100%)	1.01	25 (15%)	5 6	37, 60, 106, 144	1 (0%)
All	All	1312/1312 (100%)	0.42	108 (8%)	17 21	23, 49, 90, 144	7 (0%)

All (108) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	156	LEU	6.3
1	C	153	ILE	5.4
1	A	1	MET	4.8
1	B	156	LEU	4.8
1	G	1	MET	4.6
1	H	75	ILE	4.6
1	C	85	ILE	4.5
1	H	152	VAL	4.4
1	H	150	MET	4.3
1	G	156	LEU	4.1
1	A	154	PHE	4.0
1	B	159	TYR	4.0
1	E	75	ILE	3.8
1	D	1	MET	3.8
1	H	123	LYS	3.8
1	B	1	MET	3.7

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Mol	Chain	Res	Type	RSRZ
1	H	158	ARG	3.7
1	E	1	MET	3.6
1	E	159	TYR	3.5
1	H	108	GLU	3.4
1	H	154	PHE	3.4
1	D	153	ILE	3.3
1	F	159	TYR	3.2
1	H	1	MET	3.2
1	E	7	LYS	3.2
1	D	149	GLN	3.1
1	G	58	LYS	3.1
1	H	149	GLN	3.1
1	F	1	MET	3.1
1	G	154	PHE	3.0
1	C	159	TYR	2.9
1	H	85	ILE	2.9
1	H	86	LYS	2.9
1	A	58	LYS	2.9
1	D	80	SER	2.9
1	G	3	VAL	2.9
1	A	142	LEU	2.9
1	B	29	MET	2.9
1	H	114	LYS	2.9
1	H	160	LEU	2.8
1	G	158	ARG	2.8
1	D	154	PHE	2.8
1	C	147	ASN	2.8
1	C	1	MET	2.8
1	D	7	LYS	2.8
1	C	149	GLN	2.7
1	H	3	VAL	2.7
1	A	156	LEU	2.7
1	G	159	TYR	2.7
1	H	159	TYR	2.7
1	H	153	ILE	2.7
1	D	158	ARG	2.7
1	H	38	PHE	2.7
1	D	3	VAL	2.6
1	D	159	TYR	2.6
1	B	153	ILE	2.6
1	B	135	GLN	2.6
1	C	154	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	111	SER	2.6
1	G	7	LYS	2.6
1	E	50	GLN	2.5
1	G	155	GLN	2.5
1	B	158	ARG	2.5
1	A	3	VAL	2.5
1	H	151	SER	2.5
1	G	135	GLN	2.4
1	F	83	ASN	2.4
1	B	149	GLN	2.4
1	D	111	SER	2.4
1	F	154	PHE	2.3
1	H	41	PHE	2.3
1	B	3	VAL	2.3
1	C	123	LYS	2.3
1	F	50	GLN	2.2
1	C	144	GLU	2.2
1	F	58	LYS	2.2
1	C	141	ASP	2.2
1	G	50	GLN	2.2
1	D	83	ASN	2.2
1	G	86	LYS	2.2
1	G	75	ILE	2.2
1	G	70	VAL	2.2
1	H	7	LYS	2.2
1	H	58	LYS	2.2
1	F	153	ILE	2.2
1	C	135	GLN	2.1
1	D	155	GLN	2.1
1	G	85	ILE	2.1
1	F	3	VAL	2.1
1	H	147	ASN	2.1
1	E	156	LEU	2.1
1	A	111	SER	2.1
1	A	158	ARG	2.1
1	F	158	ARG	2.1
1	B	154	PHE	2.1
1	H	98	PHE	2.1
1	D	75	ILE	2.1
1	B	148	GLY	2.1
1	G	112	GLU	2.1
1	H	162	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	F	7	LYS	2.1
1	F	123	LYS	2.1
1	E	154	PHE	2.1
1	D	135	GLN	2.0
1	C	151	SER	2.0
1	B	39	LYS	2.0
1	G	80	SER	2.0
1	C	7	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [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	FE	D	204	1/1	0.63	0.46	280,280,280,280	0
3	GOL	G	203	6/6	0.66	0.37	93,105,124,134	0
2	SO4	A	201	5/5	0.72	0.35	142,144,146,146	0
3	GOL	C	204	6/6	0.78	0.32	86,99,119,119	0
2	SO4	H	201	5/5	0.78	0.21	72,84,106,107	5
2	SO4	H	203	5/5	0.78	0.14	99,127,135,148	0
2	SO4	C	206	5/5	0.80	0.16	107,112,121,123	5
2	SO4	B	201	5/5	0.80	0.20	146,146,150,153	0
3	GOL	E	202	6/6	0.81	0.23	55,72,97,99	0
2	SO4	D	201	5/5	0.82	0.17	95,107,113,121	0
3	GOL	G	202	6/6	0.83	0.25	57,82,100,108	0
2	SO4	B	202	5/5	0.83	0.20	69,83,108,114	0
2	SO4	B	207	5/5	0.83	0.19	114,119,120,120	5
3	GOL	A	202	6/6	0.84	0.21	55,72,105,105	0
2	SO4	H	202	5/5	0.84	0.13	65,81,92,120	5

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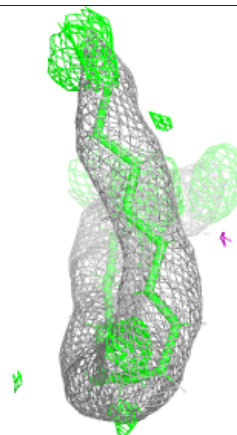
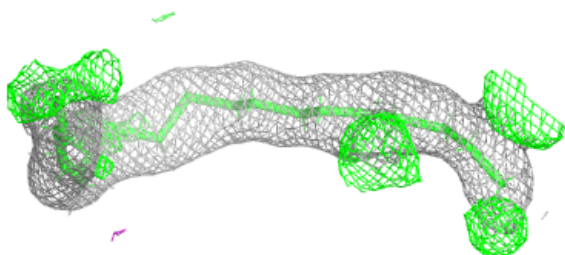
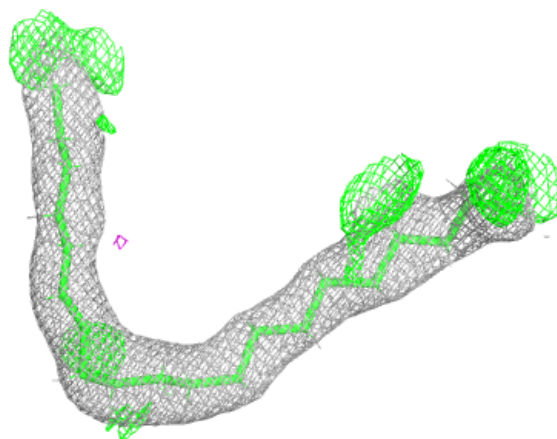
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	C	202	6/6	0.85	0.18	70,74,86,96	0
3	GOL	F	202	6/6	0.86	0.18	61,74,86,87	0
4	FE	C	201	1/1	0.86	0.15	53,53,53,53	1
3	GOL	B	205	6/6	0.86	0.19	54,71,94,94	0
2	SO4	G	205	5/5	0.88	0.20	85,85,116,117	0
2	SO4	A	204	5/5	0.89	0.15	63,65,100,117	5
2	SO4	D	203	5/5	0.89	0.14	47,56,69,77	5
3	GOL	D	202	6/6	0.89	0.15	47,56,61,66	0
5	LFA	B	206	20/20	0.89	0.21	44,69,90,100	0
5	LFA	C	205	20/20	0.89	0.19	45,67,102,107	0
3	GOL	C	203	6/6	0.90	0.16	47,67,99,119	0
5	LFA	F	203	20/20	0.91	0.20	48,69,93,116	0
5	LFA	G	204	20/20	0.91	0.21	51,80,121,123	0
2	SO4	A	203	5/5	0.92	0.15	48,51,78,85	5
3	GOL	B	203	6/6	0.93	0.12	43,52,66,67	0
2	SO4	F	201	5/5	0.93	0.10	62,69,93,94	0
2	SO4	F	204	5/5	0.94	0.10	61,62,75,81	5
4	FE	A	205	1/1	0.95	0.09	47,47,47,47	1
4	FE	G	201	1/1	0.96	0.08	55,55,55,55	1
2	SO4	B	204	5/5	0.98	0.09	44,47,54,55	5
4	FE	B	208	1/1	0.99	0.05	37,37,37,37	1
4	FE	H	204	1/1	0.99	0.11	41,41,41,41	1
4	FE	F	205	1/1	0.99	0.04	35,35,35,35	1
4	FE	E	201	1/1	1.00	0.05	30,30,30,30	1

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.

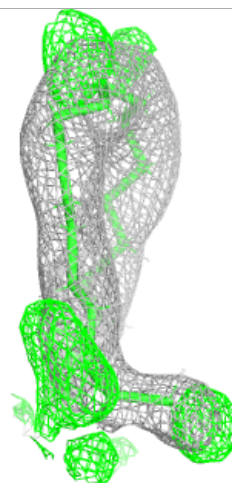
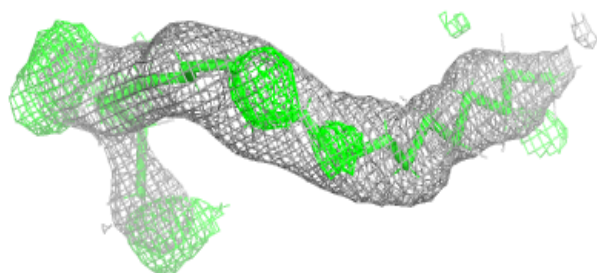
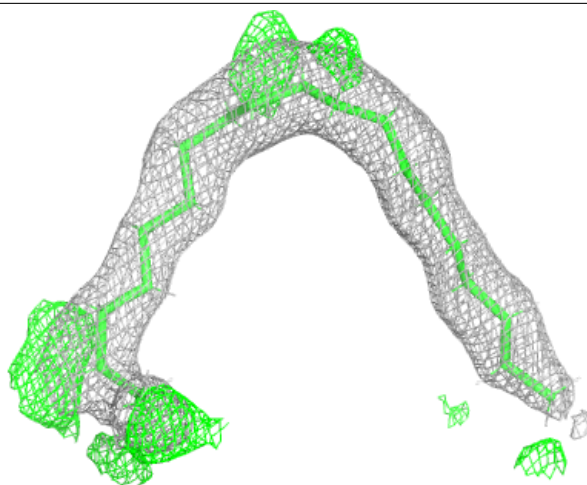
Electron density around LFA B 206:

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



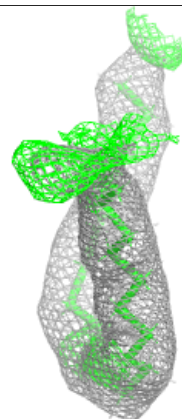
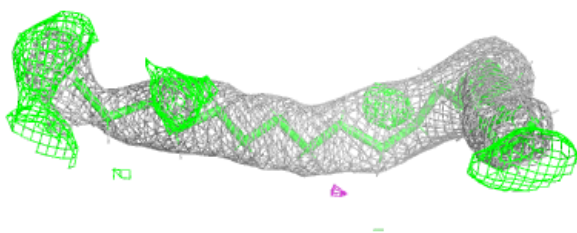
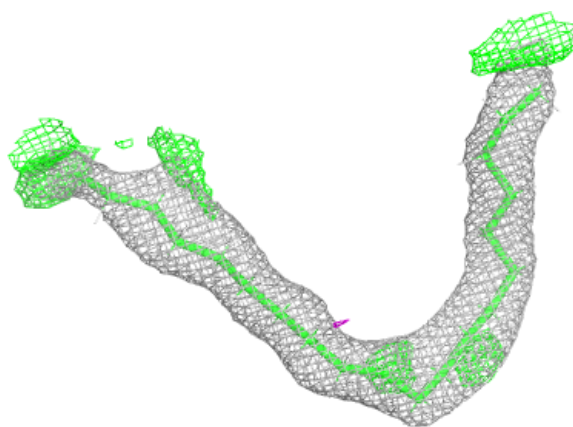
Electron density around LFA C 205:

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

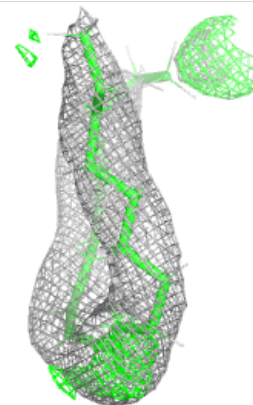
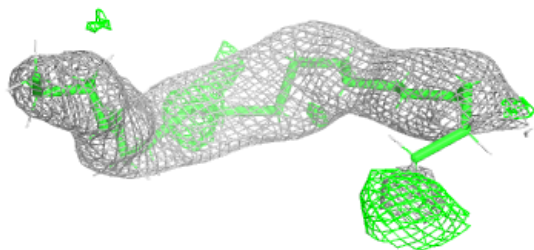
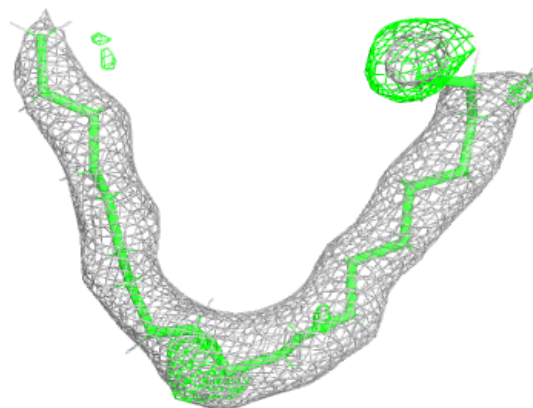


Electron density around LFA F 203:

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

**Electron density around LFA G 204:**

$2mF_o-DF_c$ (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.