Full wwPDB X-ray Structure Validation Report ⓘ

Jun 17, 2024 – 06:58 AM EDT

PDB ID	:	5OSA
Title	:	GLIC-GABAAR alpha1 chimera crystallized at pH4.6
Authors	:	Laverty, D.C.; Gold, M.G.; Smart, T.G.
Deposited on	:	2017-08-17
Resolution	:	2.75 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:

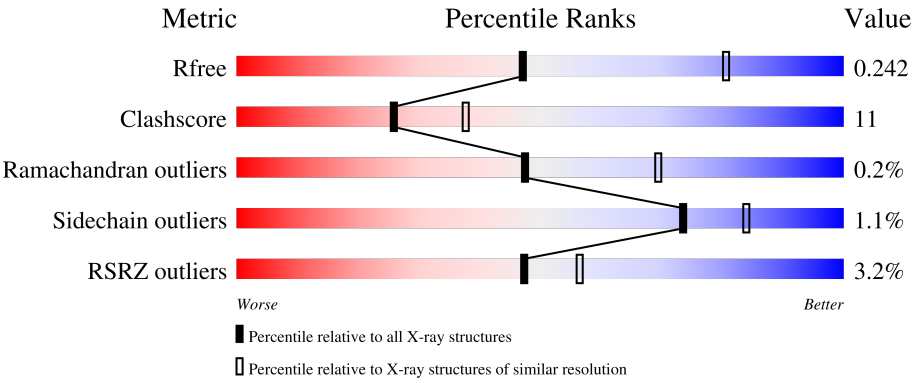
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)

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 <=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	336	4% 75% 18% 7%
1	B	336	3% 74% 18% • 7%
1	C	336	2% 75% 18% • 7%
1	D	336	4% 74% 19% • 7%
1	E	336	2% 73% 20% 7%

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	ACT	B	501	-	-	X	-
2	ACT	C	501	-	-	X	-
2	ACT	E	501	-	-	X	-
4	D12	E	504	-	-	-	X
5	Y01	D	506	-	-	-	X

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 13261 atoms, of which 459 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 Proton-gated ion channel,Gamma-aminobutyric acid receptor subunit alpha-1,Gamma-aminobutyric acid receptor subunit alpha-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	0	0
			2489	1627	400	456	6			
1	B	313	Total	C	N	O	S	0	0	0
			2495	1630	403	456	6			
1	C	313	Total	C	N	O	S	0	0	0
			2495	1630	403	456	6			
1	D	313	Total	C	N	O	S	0	0	0
			2489	1627	400	456	6			
1	E	313	Total	C	N	O	S	0	0	0
			2489	1627	400	456	6			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	258	VAL	GLY	conflict	UNP P62812
A	312	SER	-	linker	UNP P62812
A	313	GLN	-	linker	UNP P62812
A	314	PRO	-	linker	UNP P62812
A	315	ALA	-	linker	UNP P62812
A	316	ARG	-	linker	UNP P62812
A	317	ALA	-	linker	UNP P62812
A	318	ALA	-	linker	UNP P62812
A	429	HIS	-	expression tag	UNP P62812
A	430	HIS	-	expression tag	UNP P62812
A	431	HIS	-	expression tag	UNP P62812
A	432	HIS	-	expression tag	UNP P62812
A	433	HIS	-	expression tag	UNP P62812
A	434	HIS	-	expression tag	UNP P62812
A	435	HIS	-	expression tag	UNP P62812
A	436	HIS	-	expression tag	UNP P62812
B	258	VAL	GLY	conflict	UNP P62812
B	312	SER	-	linker	UNP P62812

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Chain	Residue	Modelled	Actual	Comment	Reference
B	313	GLN	-	linker	UNP P62812
B	314	PRO	-	linker	UNP P62812
B	315	ALA	-	linker	UNP P62812
B	316	ARG	-	linker	UNP P62812
B	317	ALA	-	linker	UNP P62812
B	318	ALA	-	linker	UNP P62812
B	429	HIS	-	expression tag	UNP P62812
B	430	HIS	-	expression tag	UNP P62812
B	431	HIS	-	expression tag	UNP P62812
B	432	HIS	-	expression tag	UNP P62812
B	433	HIS	-	expression tag	UNP P62812
B	434	HIS	-	expression tag	UNP P62812
B	435	HIS	-	expression tag	UNP P62812
B	436	HIS	-	expression tag	UNP P62812
C	258	VAL	GLY	conflict	UNP P62812
C	312	SER	-	linker	UNP P62812
C	313	GLN	-	linker	UNP P62812
C	314	PRO	-	linker	UNP P62812
C	315	ALA	-	linker	UNP P62812
C	316	ARG	-	linker	UNP P62812
C	317	ALA	-	linker	UNP P62812
C	318	ALA	-	linker	UNP P62812
C	429	HIS	-	expression tag	UNP P62812
C	430	HIS	-	expression tag	UNP P62812
C	431	HIS	-	expression tag	UNP P62812
C	432	HIS	-	expression tag	UNP P62812
C	433	HIS	-	expression tag	UNP P62812
C	434	HIS	-	expression tag	UNP P62812
C	435	HIS	-	expression tag	UNP P62812
C	436	HIS	-	expression tag	UNP P62812
D	258	VAL	GLY	conflict	UNP P62812
D	312	SER	-	linker	UNP P62812
D	313	GLN	-	linker	UNP P62812
D	314	PRO	-	linker	UNP P62812
D	315	ALA	-	linker	UNP P62812
D	316	ARG	-	linker	UNP P62812
D	317	ALA	-	linker	UNP P62812
D	318	ALA	-	linker	UNP P62812
D	429	HIS	-	expression tag	UNP P62812
D	430	HIS	-	expression tag	UNP P62812
D	431	HIS	-	expression tag	UNP P62812
D	432	HIS	-	expression tag	UNP P62812

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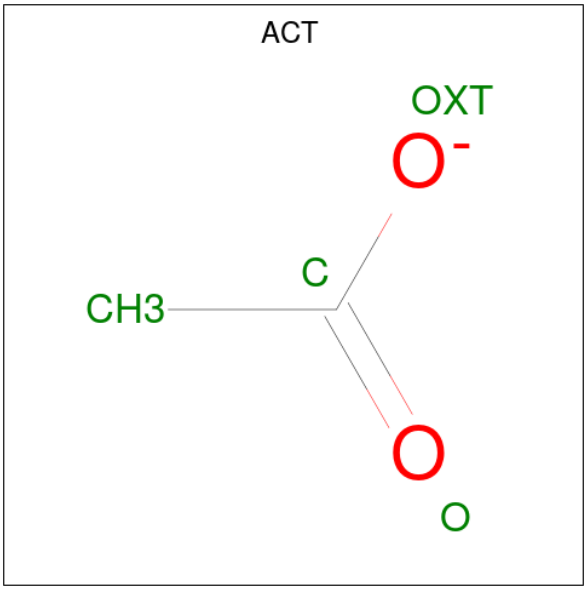
Chain	Residue	Modelled	Actual	Comment	Reference
D	433	HIS	-	expression tag	UNP P62812
D	434	HIS	-	expression tag	UNP P62812
D	435	HIS	-	expression tag	UNP P62812
D	436	HIS	-	expression tag	UNP P62812
E	258	VAL	GLY	conflict	UNP P62812
E	312	SER	-	linker	UNP P62812
E	313	GLN	-	linker	UNP P62812
E	314	PRO	-	linker	UNP P62812
E	315	ALA	-	linker	UNP P62812
E	316	ARG	-	linker	UNP P62812
E	317	ALA	-	linker	UNP P62812
E	318	ALA	-	linker	UNP P62812
E	429	HIS	-	expression tag	UNP P62812
E	430	HIS	-	expression tag	UNP P62812
E	431	HIS	-	expression tag	UNP P62812
E	432	HIS	-	expression tag	UNP P62812
E	433	HIS	-	expression tag	UNP P62812
E	434	HIS	-	expression tag	UNP P62812
E	435	HIS	-	expression tag	UNP P62812
E	436	HIS	-	expression tag	UNP P62812

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0

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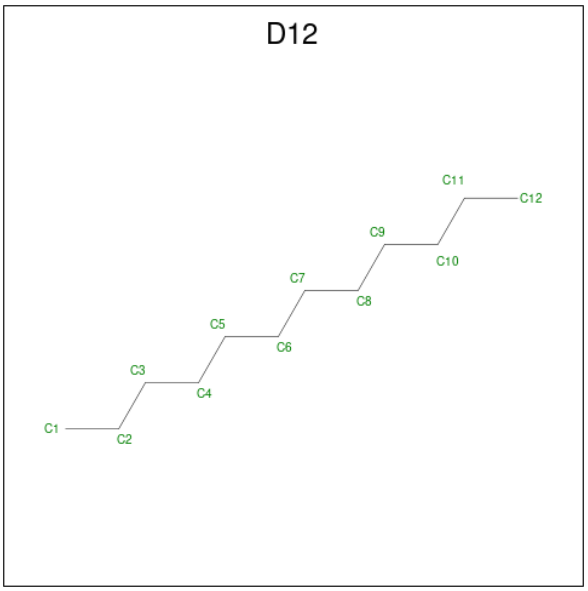
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C H O 7 2 3 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	D	1	Total C H O 7 2 3 2	0	0
2	E	1	Total C O 4 2 2	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

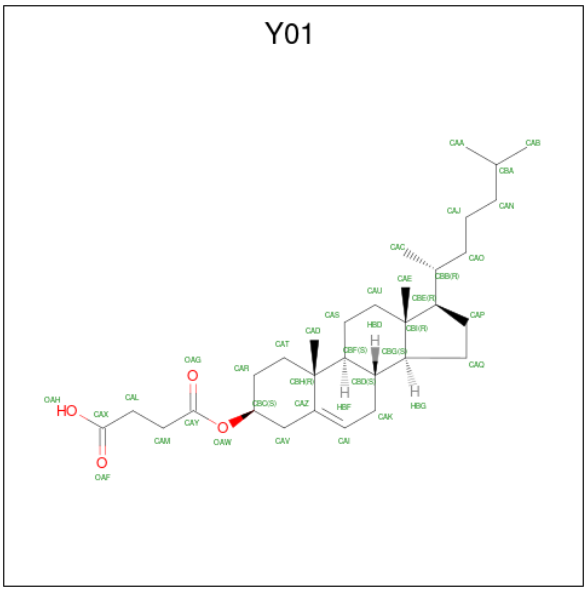
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	3	Total Cl 3 3	0	0
3	B	1	Total Cl 1 1	0	0
3	C	1	Total Cl 1 1	0	0
3	D	1	Total Cl 1 1	0	0
3	E	1	Total Cl 1 1	0	0

- Molecule 4 is DODECANE (three-letter code: D12) (formula: $C_{12}H_{26}$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C H 38 12 26	0	0
4	B	1	Total C H 38 12 26	0	0
4	B	1	Total C H 38 12 26	0	0
4	C	1	Total C H 38 12 26	0	0
4	D	1	Total C H 38 12 26	0	0
4	D	1	Total C H 38 12 26	0	0
4	E	1	Total C H 38 12 26	0	0
4	E	1	Total C H 38 12 26	0	0

- Molecule 5 is CHOLESTEROL HEMISUCCINATE (three-letter code: Y01) (formula: $C_{31}H_{50}O_4$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C H O 84 31 49 4	0	0
5	B	1	Total C H O 84 31 49 4	0	0
5	C	1	Total C H O 84 31 49 4	0	0
5	D	1	Total C H O 84 31 49 4	0	0
5	E	1	Total C H O 84 31 49 4	0	0

- Molecule 6 is water.

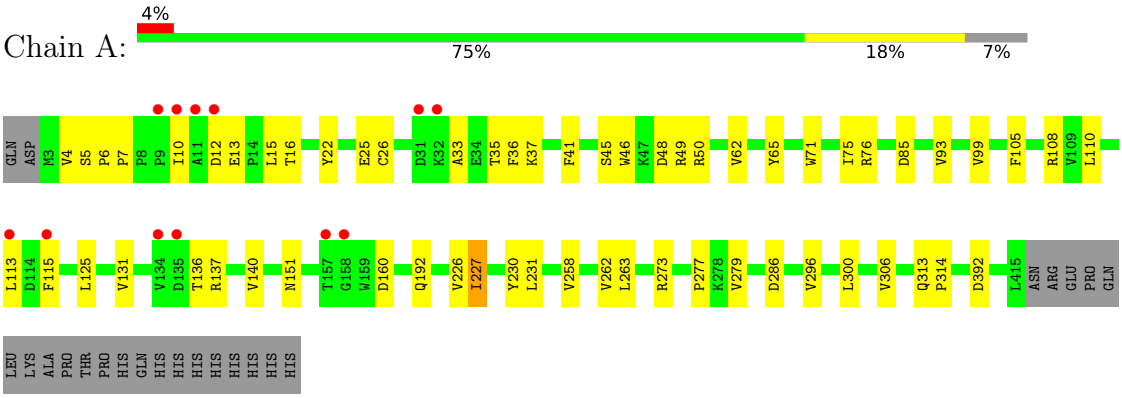
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	5	Total O 5 5	0	0
6	B	6	Total O 6 6	0	0
6	C	6	Total O 6 6	0	0
6	D	9	Total O 9 9	0	0
6	E	9	Total O 9 9	0	0

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.

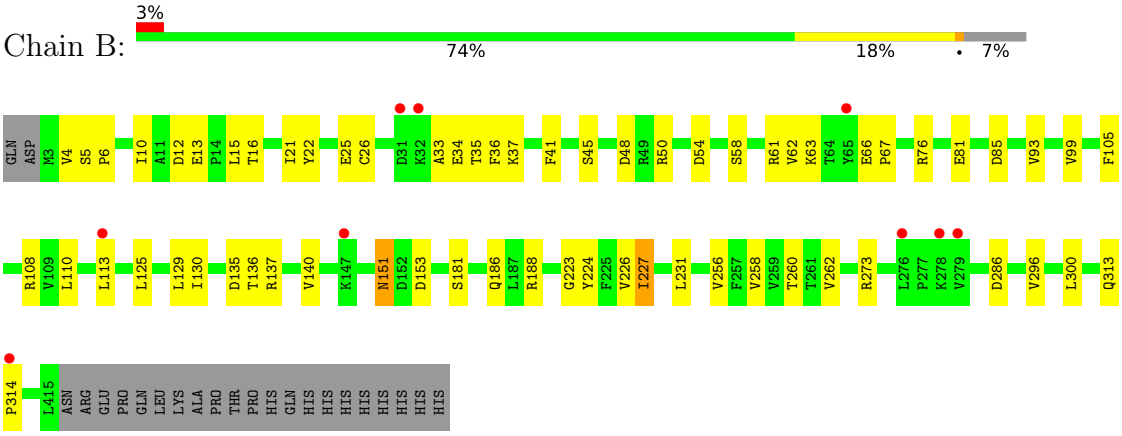
• Molecule 1: Proton-gated ion channel,Gamma-aminobutyric acid receptor subunit alpha-1,Gamma-aminobutyric acid receptor subunit alpha-1

Chain A:



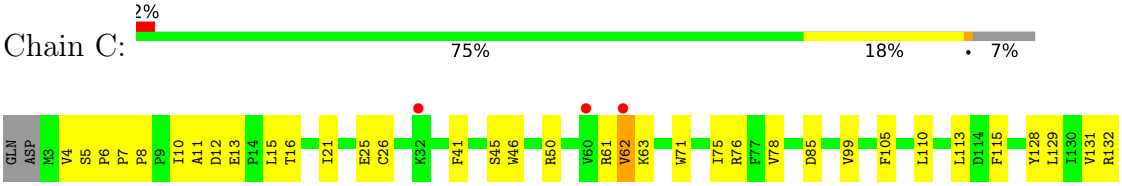
• Molecule 1: Proton-gated ion channel,Gamma-aminobutyric acid receptor subunit alpha-1,Gamma-aminobutyric acid receptor subunit alpha-1

Chain B:



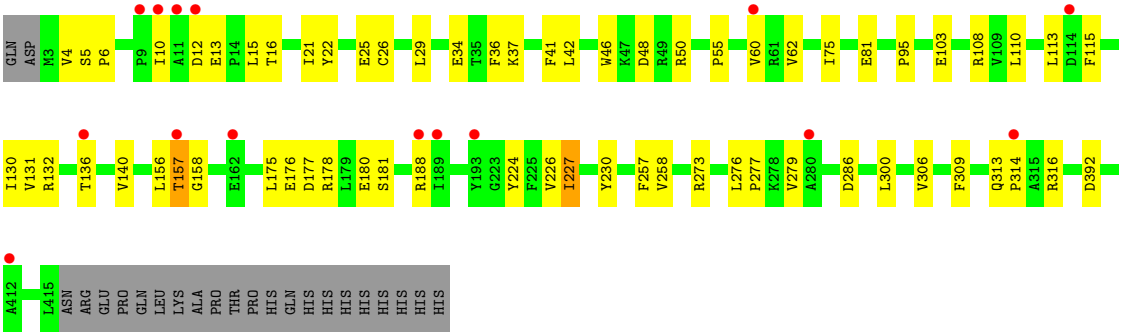
• Molecule 1: Proton-gated ion channel,Gamma-aminobutyric acid receptor subunit alpha-1,Gamma-aminobutyric acid receptor subunit alpha-1

Chain C:



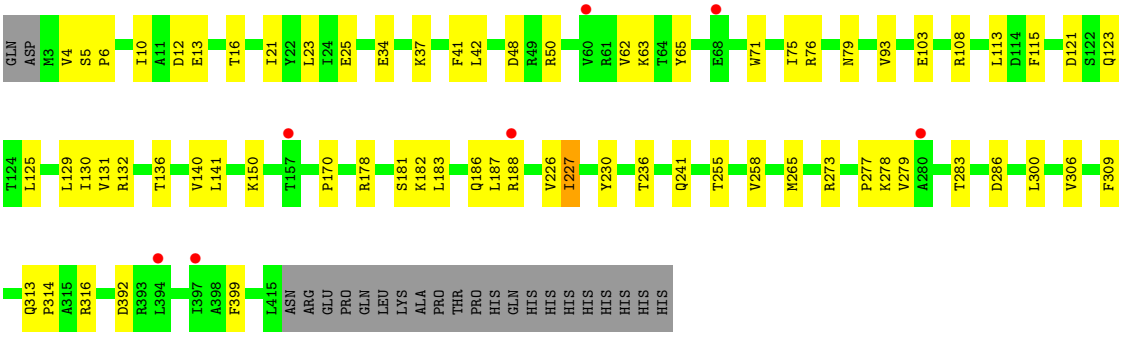
• Molecule 1: Proton-gated ion channel,Gamma-aminobutyric acid receptor subunit alpha-1,Gamma-aminobutyric acid receptor subunit alpha-1

Chain D:



• Molecule 1: Proton-gated ion channel,Gamma-aminobutyric acid receptor subunit alpha-1,Gamma-aminobutyric acid receptor subunit alpha-1

Chain E:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	185.02Å 133.95Å 162.70Å 90.00° 103.53° 90.00°	Depositor
Resolution (Å)	29.98 – 2.75 95.28 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.6 (29.98-2.75) 100.0 (95.28-2.75)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.76 (at 2.73Å)	Xtriage
Refinement program	BUSTER dev_2645, PHENIX dev_2645	Depositor
R, R_{free}	0.203 , 0.229 0.218 , 0.242	Depositor DCC
R_{free} test set	5058 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	60.7	Xtriage
Anisotropy	0.371	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 63.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.89	EDS
Total number of atoms	13261	wwPDB-VP
Average B, all atoms (Å^2)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: Y01, D12, CL, ACT

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	0.32	0/2554	0.50	0/3499
1	B	0.33	0/2560	0.50	0/3506
1	C	0.33	0/2560	0.51	0/3506
1	D	0.34	0/2554	0.52	0/3499
1	E	0.35	0/2554	0.52	0/3499
All	All	0.33	0/12782	0.51	0/17509

There are no bond length outliers.

There are no bond angle outliers.

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	2489	0	2454	53	0
1	B	2495	0	2465	57	0
1	C	2495	0	2465	58	0
1	D	2489	0	2454	59	0
1	E	2489	0	2454	67	0
2	A	8	3	6	1	0
2	B	4	0	3	3	0
2	C	8	0	6	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	8	3	6	1	0
2	E	4	0	3	2	0
3	A	3	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
4	A	12	26	26	0	0
4	B	24	52	52	0	0
4	C	12	26	26	0	0
4	D	24	52	52	0	0
4	E	24	52	52	0	0
5	A	35	49	49	3	0
5	B	35	49	49	1	0
5	C	35	49	49	5	0
5	D	35	49	49	2	0
5	E	35	49	49	1	0
6	A	5	0	0	1	0
6	B	6	0	0	0	0
6	C	6	0	0	0	0
6	D	9	0	0	1	0
6	E	9	0	0	1	0
All	All	12802	459	12769	269	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 11.

All (269) 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:E:186:GLN:OE1	1:E:188:ARG:NH2	2.01	0.93
1:A:227:ILE:HD11	1:B:273:ARG:HD2	1.56	0.86
1:D:227:ILE:HD11	1:E:273:ARG:HD2	1.56	0.86
1:A:7:PRO:O	1:A:49:ARG:NH2	2.12	0.83
1:D:37:LYS:HE3	1:D:108:ARG:HD2	1.60	0.82
1:B:186:GLN:OE1	1:B:188:ARG:NH2	2.13	0.82
1:A:273:ARG:HD2	1:E:227:ILE:HD11	1.64	0.79
1:A:110:LEU:HD23	1:E:25:GLU:HG3	1.66	0.76
1:D:157:THR:HG21	1:E:34:GLU:OE1	1.87	0.75
1:D:37:LYS:CE	1:D:108:ARG:HD2	2.17	0.74
1:D:227:ILE:HD11	1:E:273:ARG:CD	2.18	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:273:ARG:CD	1:E:227:ILE:HD11	2.18	0.73
1:B:227:ILE:HD11	1:C:273:ARG:NE	2.05	0.71
1:E:37:LYS:HE3	1:E:108:ARG:HD2	1.72	0.71
1:B:37:LYS:HE3	1:B:108:ARG:HD2	1.71	0.71
1:A:227:ILE:HD11	1:B:273:ARG:CD	2.21	0.70
1:C:157:THR:HG21	1:D:34:GLU:OE1	1.92	0.70
1:C:4:VAL:O	1:C:50:ARG:NH2	2.21	0.70
1:C:258:VAL:HG13	1:C:300:LEU:HB2	1.75	0.69
1:B:81:GLU:OE1	1:B:108:ARG:NH1	2.26	0.68
1:B:37:LYS:CE	1:B:108:ARG:HD2	2.25	0.67
5:D:506:Y01:HAC1	5:D:506:Y01:HAU2	1.77	0.67
1:A:37:LYS:HE3	1:A:108:ARG:HD2	1.77	0.67
1:E:48:ASP:OD1	1:E:50:ARG:HD3	1.96	0.65
1:A:110:LEU:HD23	1:E:25:GLU:CG	2.25	0.65
1:D:48:ASP:OD1	1:D:50:ARG:HD3	1.97	0.65
1:D:81:GLU:OE1	1:D:108:ARG:NH1	2.30	0.64
1:C:25:GLU:HG3	1:D:110:LEU:HD23	1.79	0.64
1:D:273:ARG:NH1	1:D:286:ASP:OD2	2.31	0.64
1:B:66:GLU:HG3	1:B:67:PRO:HD2	1.81	0.63
1:A:62:VAL:HG21	1:B:135:ASP:OD2	1.99	0.63
1:A:258:VAL:HG13	1:A:300:LEU:HB2	1.80	0.62
1:B:258:VAL:HG11	1:B:300:LEU:HD12	1.82	0.62
1:C:227:ILE:HD11	1:D:273:ARG:NE	2.13	0.62
1:C:227:ILE:HD11	1:D:273:ARG:CD	2.29	0.62
1:E:236:THR:HG21	1:E:265:MET:HE2	1.81	0.62
1:B:25:GLU:HG3	1:C:110:LEU:HD23	1.82	0.62
1:E:37:LYS:CE	1:E:108:ARG:HD2	2.29	0.61
1:B:151:ASN:ND2	1:B:153:ASP:OD1	2.33	0.61
1:D:6:PRO:HA	1:D:50:ARG:HH11	1.66	0.61
1:A:37:LYS:CE	1:A:108:ARG:HD2	2.30	0.61
1:C:61:ARG:HG2	1:C:62:VAL:HG22	1.82	0.60
1:D:258:VAL:HG13	1:D:300:LEU:HB2	1.84	0.60
1:A:258:VAL:HG11	1:A:300:LEU:CD1	2.31	0.60
1:B:227:ILE:HD11	1:C:273:ARG:CD	2.32	0.60
1:C:273:ARG:NH1	1:C:286:ASP:OD2	2.34	0.59
1:E:258:VAL:HG13	1:E:300:LEU:HB2	1.84	0.59
1:C:16:THR:HA	1:C:140:VAL:O	2.03	0.59
1:A:306:VAL:HG21	1:A:392:ASP:OD1	2.03	0.59
1:B:231:LEU:HD13	5:C:505:Y01:HAR1	1.85	0.59
1:A:48:ASP:OD1	1:A:50:ARG:HD3	2.03	0.59
1:C:156:LEU:O	1:C:158:GLY:N	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:6:PRO:HA	1:B:50:ARG:HH11	1.68	0.58
1:B:48:ASP:OD1	1:B:50:ARG:HD3	2.03	0.58
1:D:16:THR:HA	1:D:140:VAL:O	2.03	0.58
1:B:62:VAL:HG23	1:B:63:LYS:N	2.19	0.58
1:E:65:TYR:HE1	1:E:93:VAL:HG21	1.69	0.58
5:C:505:Y01:HAC1	5:C:505:Y01:HAE2	1.84	0.58
1:D:55:PRO:HG3	1:D:95:PRO:CG	2.33	0.58
1:B:10:ILE:HG22	1:B:10:ILE:O	2.03	0.58
1:E:258:VAL:HG11	1:E:300:LEU:CD1	2.34	0.58
1:A:33:ALA:O	1:A:35:THR:HG23	2.04	0.58
1:E:273:ARG:NH1	1:E:286:ASP:OD2	2.38	0.57
1:C:25:GLU:CG	1:D:110:LEU:HD23	2.35	0.57
1:E:62:VAL:HG23	1:E:63:LYS:N	2.19	0.56
1:C:76:ARG:NH1	2:C:501:ACT:O	2.39	0.56
1:A:258:VAL:HG11	1:A:300:LEU:HD12	1.87	0.56
1:E:76:ARG:NH1	2:E:501:ACT:O	2.38	0.56
1:D:157:THR:HG22	1:D:157:THR:O	2.05	0.56
1:D:227:ILE:CD1	1:E:273:ARG:HD2	2.31	0.56
1:A:36:PHE:CE1	1:A:125:LEU:HB3	2.40	0.56
1:D:156:LEU:O	1:D:158:GLY:N	2.38	0.56
1:B:227:ILE:HD11	1:C:273:ARG:CZ	2.34	0.56
1:B:262:VAL:HG22	1:B:296:VAL:HG12	1.88	0.56
1:D:75:ILE:HD13	1:D:131:VAL:HB	1.87	0.56
1:D:258:VAL:HG11	1:D:300:LEU:HD12	1.88	0.55
1:B:61:ARG:HG3	1:B:62:VAL:HG13	1.87	0.55
1:B:34:GLU:HG2	1:B:113:LEU:HG	1.88	0.55
1:C:15:LEU:HD11	1:C:46:TRP:HB2	1.88	0.55
1:D:60:VAL:HG12	1:D:62:VAL:H	1.71	0.55
1:E:258:VAL:HG11	1:E:300:LEU:HD12	1.87	0.55
1:B:273:ARG:NH1	1:B:286:ASP:OD2	2.40	0.55
1:A:16:THR:HA	1:A:140:VAL:O	2.06	0.54
1:E:10:ILE:HB	1:E:13:GLU:HB2	1.87	0.54
1:B:258:VAL:HG13	1:B:300:LEU:HB2	1.89	0.54
1:E:76:ARG:NH2	6:E:601:HOH:O	2.39	0.54
1:E:226:VAL:HA	1:E:230:TYR:HB2	1.90	0.54
1:B:256:VAL:O	1:B:260:THR:HG23	2.07	0.54
1:C:8:PRO:HG2	1:C:11:ALA:HA	1.90	0.54
1:C:134:VAL:HG23	1:C:137:ARG:H	1.71	0.54
1:D:130:ILE:HD13	2:D:501:ACT:H2	1.88	0.54
1:A:15:LEU:HD11	1:A:46:TRP:HB2	1.89	0.54
1:D:10:ILE:HG22	1:D:10:ILE:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:PRO:HG2	1:A:15:LEU:HD22	1.89	0.53
1:B:10:ILE:HB	1:B:13:GLU:HB2	1.89	0.53
1:A:6:PRO:HA	1:A:50:ARG:HH11	1.72	0.53
1:A:10:ILE:HB	1:A:13:GLU:HB2	1.89	0.53
1:A:10:ILE:O	1:A:10:ILE:HG22	2.07	0.53
1:E:75:ILE:HD13	1:E:131:VAL:HB	1.91	0.53
1:D:175:LEU:HB3	1:D:180:GLU:HG3	1.91	0.53
1:A:273:ARG:NE	1:E:227:ILE:HD11	2.23	0.53
1:C:258:VAL:HG11	1:C:300:LEU:CD1	2.39	0.53
1:E:34:GLU:HG2	1:E:113:LEU:HG	1.90	0.52
1:E:16:THR:HA	1:E:140:VAL:O	2.10	0.52
1:C:10:ILE:O	1:C:10:ILE:HG22	2.08	0.52
1:B:6:PRO:HA	1:B:50:ARG:NH1	2.25	0.52
1:C:227:ILE:HD11	1:D:273:ARG:HD2	1.90	0.52
1:D:34:GLU:HG2	1:D:113:LEU:HG	1.91	0.52
1:E:313:GLN:N	1:E:314:PRO:HD3	2.24	0.52
1:A:6:PRO:HA	1:A:50:ARG:NH1	2.24	0.52
1:E:6:PRO:HA	1:E:50:ARG:HH11	1.76	0.51
1:A:258:VAL:CG1	1:A:300:LEU:HD13	2.40	0.51
1:C:262:VAL:HG22	1:C:296:VAL:HG12	1.92	0.51
1:C:6:PRO:HA	1:C:50:ARG:HH11	1.74	0.51
1:D:6:PRO:HA	1:D:50:ARG:NH1	2.24	0.51
5:A:507:Y01:HAU2	5:A:507:Y01:HAC1	1.93	0.51
1:B:25:GLU:CG	1:C:110:LEU:HD23	2.40	0.51
1:D:277:PRO:HG2	1:D:279:VAL:HG13	1.93	0.51
1:B:224:TYR:CD2	1:C:278:LYS:HA	2.47	0.51
1:D:306:VAL:HG21	1:D:392:ASP:OD1	2.10	0.50
1:C:258:VAL:CG1	1:C:300:LEU:HD13	2.41	0.50
1:A:4:VAL:O	1:A:50:ARG:NH2	2.40	0.50
1:C:258:VAL:HG11	1:C:300:LEU:HD12	1.94	0.50
1:E:4:VAL:HG22	1:E:5:SER:H	1.76	0.50
1:C:45:SER:HA	1:C:99:VAL:O	2.12	0.50
1:E:227:ILE:HG13	1:E:227:ILE:O	2.11	0.50
1:A:227:ILE:CD1	1:B:273:ARG:HD2	2.37	0.50
1:E:258:VAL:CG1	1:E:300:LEU:HD13	2.42	0.50
1:A:65:TYR:HE1	1:A:93:VAL:HG21	1.77	0.49
1:B:15:LEU:HB2	1:B:137:ARG:CZ	2.43	0.49
1:C:313:GLN:N	1:C:314:PRO:HD3	2.27	0.49
1:E:76:ARG:HH11	2:E:501:ACT:C	2.25	0.49
1:A:76:ARG:NH1	2:A:501:ACT:O	2.46	0.49
5:B:505:Y01:HAC1	5:B:505:Y01:HAU2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:THR:HA	1:B:140:VAL:O	2.13	0.49
1:D:227:ILE:HD11	1:E:273:ARG:NE	2.27	0.49
1:B:258:VAL:HG11	1:B:300:LEU:CD1	2.43	0.48
1:D:4:VAL:HG22	1:D:5:SER:H	1.78	0.48
1:C:6:PRO:HA	1:C:50:ARG:NH1	2.28	0.48
1:A:258:VAL:CG1	1:A:300:LEU:HB2	2.43	0.48
1:A:273:ARG:NH1	1:A:286:ASP:OD2	2.46	0.48
1:A:226:VAL:HA	1:A:230:TYR:HB2	1.95	0.48
1:C:10:ILE:HB	1:C:13:GLU:HB2	1.94	0.48
1:D:37:LYS:HE2	1:D:108:ARG:HD2	1.96	0.48
1:D:15:LEU:HD11	1:D:46:TRP:HB2	1.95	0.48
1:B:76:ARG:NH1	2:B:501:ACT:O	2.48	0.47
1:A:110:LEU:CD2	1:E:25:GLU:HG3	2.40	0.47
1:A:45:SER:HA	1:A:99:VAL:O	2.15	0.47
1:A:231:LEU:HA	1:A:231:LEU:HD23	1.69	0.47
1:E:129:LEU:O	1:E:182:LYS:HA	2.14	0.47
1:A:151:ASN:ND2	6:A:602:HOH:O	2.48	0.47
1:D:10:ILE:HB	1:D:13:GLU:HB2	1.96	0.47
1:D:313:GLN:N	1:D:314:PRO:HD3	2.30	0.47
1:E:236:THR:CG2	1:E:265:MET:HE2	2.43	0.47
1:A:75:ILE:HD13	1:A:131:VAL:HB	1.97	0.47
1:A:25:GLU:HG3	1:B:110:LEU:HD23	1.97	0.46
1:B:258:VAL:CG1	1:B:300:LEU:CD1	2.93	0.46
1:C:156:LEU:C	1:C:158:GLY:H	2.18	0.46
1:E:10:ILE:HG22	1:E:10:ILE:O	2.15	0.46
1:E:42:LEU:HB3	1:E:103:GLU:HG2	1.98	0.46
1:C:75:ILE:HD13	1:C:131:VAL:HB	1.96	0.46
1:D:309:PHE:CD2	1:D:316:ARG:HD2	2.51	0.46
1:A:160:ASP:OD1	1:A:192:GLN:NE2	2.45	0.46
1:A:262:VAL:HG22	1:A:296:VAL:HG12	1.97	0.46
1:B:231:LEU:HD23	1:B:231:LEU:HA	1.74	0.46
1:C:10:ILE:C	1:C:12:ASP:H	2.18	0.46
1:E:236:THR:CG2	1:E:265:MET:CE	2.92	0.46
1:A:258:VAL:CG1	1:A:300:LEU:CD1	2.94	0.46
1:D:55:PRO:HG3	1:D:95:PRO:HG3	1.98	0.46
1:D:258:VAL:CG1	1:D:300:LEU:CD1	2.94	0.46
1:D:75:ILE:CD1	1:D:131:VAL:HB	2.46	0.46
1:C:113:LEU:HB2	1:C:115:PHE:CE2	2.50	0.46
1:A:85:ASP:O	1:A:105:PHE:HA	2.16	0.45
1:B:76:ARG:HH11	2:B:501:ACT:C	2.28	0.45
1:C:7:PRO:HD2	1:C:50:ARG:HD2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:10:ILE:C	1:D:12:ASP:H	2.19	0.45
1:E:10:ILE:HD12	1:E:13:GLU:CD	2.36	0.45
1:B:62:VAL:CG2	1:B:63:LYS:N	2.80	0.45
1:B:227:ILE:HD11	1:C:273:ARG:HD2	1.98	0.45
1:B:85:ASP:O	1:B:105:PHE:HA	2.17	0.45
1:B:258:VAL:HG11	1:B:300:LEU:HA	1.98	0.45
1:A:313:GLN:N	1:A:314:PRO:HD3	2.32	0.45
1:C:258:VAL:CG1	1:C:300:LEU:HB2	2.46	0.45
1:E:125:LEU:HD12	1:E:187:LEU:HD23	1.98	0.45
5:E:505:Y01:HAC1	5:E:505:Y01:HAU2	1.99	0.45
1:D:25:GLU:OE1	1:E:79:ASN:HA	2.17	0.44
1:E:258:VAL:CG1	1:E:300:LEU:CD1	2.95	0.44
1:B:37:LYS:HE2	1:B:108:ARG:HD2	1.99	0.44
1:C:76:ARG:HH11	2:C:501:ACT:C	2.31	0.44
1:C:258:VAL:CG1	1:C:300:LEU:CD1	2.96	0.44
1:E:75:ILE:CD1	1:E:131:VAL:HB	2.47	0.44
1:B:130:ILE:HA	1:B:181:SER:O	2.18	0.44
1:C:227:ILE:HD11	1:D:273:ARG:CZ	2.48	0.44
1:D:42:LEU:HB3	1:D:103:GLU:HG2	2.00	0.44
1:D:258:VAL:HG11	1:D:300:LEU:CD1	2.47	0.44
1:A:15:LEU:HB2	1:A:137:ARG:CZ	2.48	0.44
1:B:4:VAL:HG22	1:B:5:SER:H	1.83	0.44
1:B:45:SER:HA	1:B:99:VAL:O	2.18	0.43
1:E:309:PHE:CD2	1:E:316:ARG:HD2	2.53	0.43
1:B:21:ILE:HA	1:B:41:PHE:O	2.18	0.43
1:D:21:ILE:HA	1:D:41:PHE:O	2.17	0.43
1:E:10:ILE:C	1:E:12:ASP:H	2.22	0.43
1:E:130:ILE:HA	1:E:181:SER:O	2.18	0.43
1:B:10:ILE:C	1:B:12:ASP:H	2.21	0.43
1:C:277:PRO:HG2	1:C:279:VAL:HG13	1.99	0.43
1:D:227:ILE:HD11	1:E:273:ARG:CZ	2.49	0.43
1:E:115:PHE:O	1:E:283:THR:HG22	2.17	0.43
1:B:76:ARG:HD2	2:B:501:ACT:OXT	2.17	0.43
1:C:4:VAL:HG22	1:C:5:SER:H	1.82	0.43
1:C:157:THR:O	1:C:157:THR:HG22	2.18	0.43
1:E:129:LEU:HD23	1:E:129:LEU:HA	1.83	0.43
1:E:121:ASP:HB2	1:E:123:GLN:NE2	2.34	0.43
1:A:113:LEU:HB2	1:A:115:PHE:CE2	2.54	0.43
1:C:132:ARG:CZ	1:C:178:ARG:HB2	2.48	0.43
1:D:257:PHE:CE1	1:E:255:THR:HG23	2.54	0.43
1:B:36:PHE:CE1	1:B:125:LEU:HB3	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:241:GLN:NE2	1:E:399:PHE:HB3	2.33	0.42
1:A:22:TYR:HB3	1:A:41:PHE:HB2	2.00	0.42
1:B:93:VAL:HG12	1:B:99:VAL:HG22	2.01	0.42
1:C:290:ALA:HA	5:C:505:Y01:CAM	2.49	0.42
1:D:22:TYR:HB3	1:D:41:PHE:HB2	2.00	0.42
1:E:42:LEU:HB3	1:E:103:GLU:CG	2.49	0.42
5:A:507:Y01:HAC1	5:A:507:Y01:HAE2	1.99	0.42
1:C:21:ILE:HA	1:C:41:PHE:O	2.18	0.42
1:D:224:TYR:CD2	1:E:278:LYS:HA	2.54	0.42
1:C:6:PRO:HD3	1:C:71:TRP:CE3	2.54	0.42
1:C:129:LEU:O	1:C:182:LYS:HA	2.20	0.42
1:E:6:PRO:HD3	1:E:71:TRP:CE3	2.55	0.42
1:E:132:ARG:CZ	1:E:178:ARG:HB2	2.50	0.42
5:C:505:Y01:HAC1	5:C:505:Y01:HAU2	2.02	0.42
1:E:141:LEU:O	1:E:170:PRO:HG3	2.19	0.42
1:A:6:PRO:HD3	1:A:71:TRP:CE3	2.55	0.42
1:E:258:VAL:CG1	1:E:300:LEU:HB2	2.50	0.41
1:B:129:LEU:HD23	1:B:129:LEU:HA	1.85	0.41
1:C:156:LEU:HD23	1:C:156:LEU:HA	1.82	0.41
1:C:231:LEU:HA	1:C:231:LEU:HD23	1.77	0.41
1:D:29:LEU:HD13	1:D:36:PHE:HB3	2.02	0.41
5:A:507:Y01:CAP	5:A:507:Y01:CAJ	2.98	0.41
1:B:223:GLY:O	1:B:226:VAL:HB	2.20	0.41
1:C:141:LEU:O	1:C:170:PRO:HG3	2.21	0.41
1:C:290:ALA:HB2	5:C:505:Y01:HAM1	2.02	0.41
1:E:121:ASP:OD2	1:E:123:GLN:NE2	2.51	0.41
1:A:65:TYR:HE1	1:A:93:VAL:CG2	2.33	0.41
1:B:22:TYR:HB3	1:B:41:PHE:HB2	2.03	0.41
1:D:132:ARG:CZ	1:D:178:ARG:HB2	2.50	0.41
1:D:176:GLU:O	1:D:177:ASP:HB2	2.21	0.41
1:A:4:VAL:HG22	1:A:5:SER:H	1.85	0.41
1:A:263:LEU:HD23	1:A:263:LEU:HA	1.86	0.41
1:B:33:ALA:O	1:B:35:THR:HG23	2.21	0.41
1:C:398:ALA:O	1:C:402:LEU:HD23	2.21	0.41
1:D:130:ILE:HA	1:D:181:SER:O	2.21	0.41
1:B:313:GLN:N	1:B:314:PRO:HD3	2.36	0.41
1:C:78:VAL:HB	1:C:128:TYR:HB2	2.02	0.41
1:D:188:ARG:NH1	6:D:602:HOH:O	2.48	0.41
1:E:23:LEU:HB2	1:E:150:LYS:CB	2.50	0.41
5:D:506:Y01:HAC1	5:D:506:Y01:CAU	2.48	0.41
1:B:54:ASP:O	1:B:58:SER:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:62:VAL:HG23	1:E:63:LYS:H	1.83	0.40
1:D:226:VAL:HA	1:D:230:TYR:HB2	2.03	0.40
1:D:258:VAL:CG1	1:D:300:LEU:HD13	2.51	0.40
1:A:10:ILE:C	1:A:12:ASP:H	2.24	0.40
1:C:115:PHE:O	1:C:283:THR:HG22	2.20	0.40
1:E:306:VAL:HG21	1:E:392:ASP:OD1	2.20	0.40
1:A:277:PRO:HG2	1:A:279:VAL:HG13	2.03	0.40
1:D:115:PHE:CD1	1:D:276:LEU:HD22	2.57	0.40
1:E:183:LEU:HD23	1:E:183:LEU:HA	1.92	0.40
1:E:277:PRO:HG2	1:E:279:VAL:HG13	2.03	0.40
1:C:85:ASP:O	1:C:105:PHE:HA	2.21	0.40
1:E:21:ILE:HA	1:E:41:PHE:O	2.21	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	311/336 (93%)	298 (96%)	13 (4%)	0	100	100
1	B	311/336 (93%)	299 (96%)	12 (4%)	0	100	100
1	C	311/336 (93%)	298 (96%)	11 (4%)	2 (1%)	25	42
1	D	311/336 (93%)	299 (96%)	11 (4%)	1 (0%)	41	60
1	E	311/336 (93%)	299 (96%)	12 (4%)	0	100	100
All	All	1555/1680 (93%)	1493 (96%)	59 (4%)	3 (0%)	47	69

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	157	THR
1	D	157	THR

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Mol	Chain	Res	Type
1	C	63	LYS

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	271/302 (90%)	268 (99%)	3 (1%)	73	84
1	B	272/302 (90%)	268 (98%)	4 (2%)	65	78
1	C	272/302 (90%)	269 (99%)	3 (1%)	73	84
1	D	271/302 (90%)	268 (99%)	3 (1%)	73	84
1	E	271/302 (90%)	269 (99%)	2 (1%)	84	89
All	All	1357/1510 (90%)	1342 (99%)	15 (1%)	73	84

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	CYS
1	A	136	THR
1	A	227	ILE
1	B	26	CYS
1	B	136	THR
1	B	151	ASN
1	B	227	ILE
1	C	26	CYS
1	C	62	VAL
1	C	227	ILE
1	D	26	CYS
1	D	136	THR
1	D	227	ILE
1	E	136	THR
1	E	227	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	151	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry

Of 28 ligands modelled in this entry, 7 are monoatomic - leaving 21 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	ACT	C	502	-	3,3,3	1.47	1 (33%)	3,3,3	1.49	0
4	D12	D	505	-	11,11,11	0.36	0	10,10,10	0.68	0
2	ACT	D	502	-	3,3,3	1.42	1 (33%)	3,3,3	1.51	0
2	ACT	D	501	-	3,3,3	1.40	1 (33%)	3,3,3	1.56	1 (33%)
4	D12	B	503	-	11,11,11	0.38	0	10,10,10	0.68	0
5	Y01	A	507	-	38,38,38	0.61	0	57,57,57	1.34	9 (15%)
5	Y01	C	505	-	38,38,38	0.58	0	57,57,57	1.37	9 (15%)
2	ACT	A	502	-	3,3,3	1.43	1 (33%)	3,3,3	1.35	0
5	Y01	E	505	-	38,38,38	0.61	0	57,57,57	1.32	7 (12%)
2	ACT	B	501	-	3,3,3	1.38	1 (33%)	3,3,3	1.49	0
4	D12	A	506	-	11,11,11	0.41	0	10,10,10	0.64	0
4	D12	D	504	-	11,11,11	0.38	0	10,10,10	0.70	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
5	Y01	B	505	-	38,38,38	0.61	0	57,57,57	1.36	8 (14%)
2	ACT	A	501	-	3,3,3	1.43	1 (33%)	3,3,3	1.36	0
2	ACT	C	501	-	3,3,3	1.29	0	3,3,3	1.37	0
4	D12	E	504	-	11,11,11	0.36	0	10,10,10	0.72	0
4	D12	B	504	-	11,11,11	0.38	0	10,10,10	0.71	0
5	Y01	D	506	-	38,38,38	0.60	0	57,57,57	1.29	9 (15%)
4	D12	E	503	-	11,11,11	0.37	0	10,10,10	0.75	0
2	ACT	E	501	-	3,3,3	1.47	1 (33%)	3,3,3	1.35	0
4	D12	C	504	-	11,11,11	0.37	0	10,10,10	0.71	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
4	D12	E	504	-	-	1/9/9/9	-
5	Y01	A	507	-	-	3/19/77/77	0/4/4/4
5	Y01	C	505	-	-	9/19/77/77	0/4/4/4
5	Y01	E	505	-	-	3/19/77/77	0/4/4/4
4	D12	B	504	-	-	0/9/9/9	-
5	Y01	D	506	-	-	1/19/77/77	0/4/4/4
4	D12	A	506	-	-	1/9/9/9	-
4	D12	E	503	-	-	1/9/9/9	-
4	D12	D	504	-	-	0/9/9/9	-
4	D12	C	504	-	-	2/9/9/9	-
5	Y01	B	505	-	-	3/19/77/77	0/4/4/4
4	D12	B	503	-	-	1/9/9/9	-
4	D12	D	505	-	-	2/9/9/9	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	502	ACT	CH3-C	2.28	1.58	1.49
2	A	502	ACT	CH3-C	2.20	1.58	1.49
2	D	502	ACT	CH3-C	2.17	1.58	1.49
2	A	501	ACT	CH3-C	2.12	1.58	1.49
2	D	501	ACT	CH3-C	2.10	1.57	1.49
2	E	501	ACT	CH3-C	2.10	1.57	1.49
2	B	501	ACT	CH3-C	2.03	1.57	1.49

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	C	505	Y01	OAW-CAY-CAM	3.75	119.59	111.50
5	B	505	Y01	OAW-CAY-CAM	3.40	118.83	111.50
5	A	507	Y01	OAW-CAY-CAM	3.09	118.16	111.50
5	E	505	Y01	OAW-CAY-CAM	3.07	118.11	111.50
5	D	506	Y01	CAC-CBB-CBE	-2.98	108.36	112.92
5	A	507	Y01	CAQ-CBG-CBD	-2.91	114.29	119.08
5	B	505	Y01	CAK-CBD-CBG	-2.85	106.78	110.91
5	C	505	Y01	CAC-CBB-CAO	-2.82	105.95	110.36
5	C	505	Y01	CAQ-CBG-CBI	-2.74	100.54	103.84
5	D	506	Y01	CBI-CBE-CBB	-2.72	115.23	119.49
5	D	506	Y01	CAM-CAL-CAX	-2.70	107.80	113.60
5	E	505	Y01	CAQ-CBG-CBI	-2.66	100.64	103.84
5	E	505	Y01	CAK-CBD-CBG	-2.65	107.06	110.91
5	A	507	Y01	CBI-CBE-CBB	-2.64	115.34	119.49
5	A	507	Y01	CAM-CAL-CAX	-2.64	107.92	113.60
5	D	506	Y01	OAW-CAY-CAM	2.63	117.17	111.50
5	D	506	Y01	CAQ-CBG-CBI	-2.62	100.68	103.84
5	A	507	Y01	CAC-CBB-CAO	-2.55	106.36	110.36
5	A	507	Y01	CAV-CAZ-CAI	-2.52	116.98	120.61
5	B	505	Y01	CAV-CAZ-CAI	-2.51	116.99	120.61
5	C	505	Y01	CBC-CAV-CAZ	-2.51	107.62	111.52
5	E	505	Y01	CBG-CBI-CBE	-2.44	97.18	100.07
5	C	505	Y01	CBI-CBE-CBB	-2.39	115.74	119.49
5	E	505	Y01	CAM-CAL-CAX	-2.37	108.51	113.60
5	C	505	Y01	CAS-CAU-CBI	-2.35	108.75	112.78
5	C	505	Y01	CAQ-CBG-CBD	-2.32	115.26	119.08
5	D	506	Y01	CAK-CBD-CBG	-2.31	107.56	110.91
5	B	505	Y01	CAQ-CBG-CBD	-2.30	115.29	119.08
5	B	505	Y01	CBI-CBE-CBB	-2.27	115.94	119.49
5	A	507	Y01	CAK-CBD-CBG	-2.17	107.77	110.91
5	A	507	Y01	CAP-CAQ-CBG	-2.17	100.84	105.13
5	B	505	Y01	CAV-CAZ-CBH	2.15	119.28	116.42
5	A	507	Y01	CAQ-CBG-CBI	-2.14	101.27	103.84
5	D	506	Y01	CAV-CAZ-CAI	-2.13	117.53	120.61
5	B	505	Y01	CBG-CBI-CBE	-2.11	97.57	100.07
5	C	505	Y01	CBF-CBH-CAZ	2.09	112.92	109.65
5	E	505	Y01	OAH-CAX-CAL	2.08	120.71	114.03
5	D	506	Y01	OAH-CAX-CAL	2.04	120.59	114.03
5	D	506	Y01	CBC-OAW-CAY	2.03	122.79	117.79
5	C	505	Y01	CBD-CAK-CAI	-2.03	109.82	112.73
5	E	505	Y01	CBF-CBH-CAZ	2.02	112.82	109.65
5	B	505	Y01	OAH-CAX-CAL	2.01	120.50	114.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	501	ACT	O-C-CH3	-2.00	114.54	122.33

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	505	Y01	CAJ-CAO-CBB-CBE
5	B	505	Y01	CAJ-CAO-CBB-CAC
5	C	505	Y01	CAO-CBB-CBE-CAP
5	E	505	Y01	CAO-CAJ-CAN-CBA
5	E	505	Y01	CAN-CAJ-CAO-CBB
5	C	505	Y01	CAX-CAL-CAM-CAY
5	D	506	Y01	CAN-CAJ-CAO-CBB
4	A	506	D12	C11-C10-C9-C8
4	C	504	D12	C6-C7-C8-C9
5	C	505	Y01	CAM-CAY-OAW-CBC
5	A	507	Y01	CAO-CAJ-CAN-CBA
5	C	505	Y01	OAG-CAY-OAW-CBC
4	D	505	D12	C5-C6-C7-C8
5	E	505	Y01	CAO-CBB-CBE-CAP
5	B	505	Y01	CAN-CAJ-CAO-CBB
5	C	505	Y01	CAC-CBB-CBE-CAP
5	A	507	Y01	CAO-CBB-CBE-CAP
4	E	503	D12	C3-C4-C5-C6
5	A	507	Y01	CAJ-CAO-CBB-CAC
5	C	505	Y01	CAJ-CAO-CBB-CAC
4	E	504	D12	C11-C10-C9-C8
5	C	505	Y01	CAM-CAL-CAX-OAH
4	C	504	D12	C1-C2-C3-C4
4	B	503	D12	C2-C3-C4-C5
5	C	505	Y01	CAM-CAL-CAX-OAF
4	D	505	D12	C7-C8-C9-C10
5	C	505	Y01	CAO-CAJ-CAN-CBA

There are no ring outliers.

10 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	ACT	1	0
5	A	507	Y01	3	0
5	C	505	Y01	5	0

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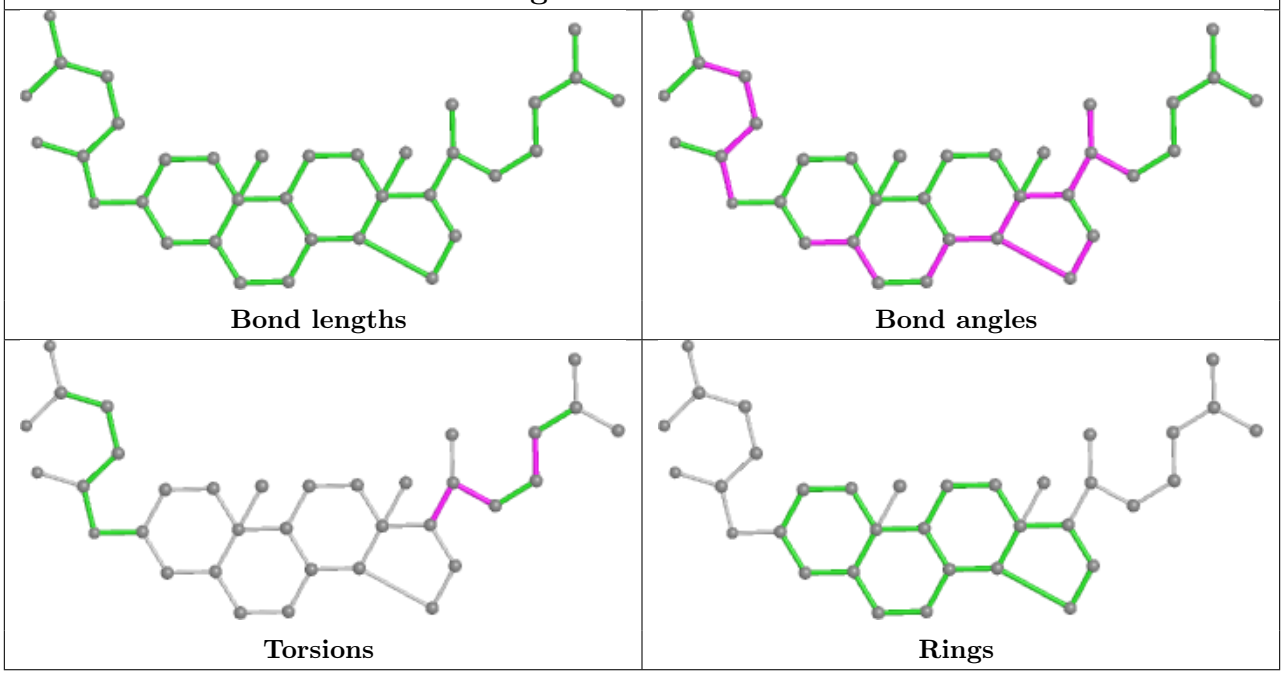
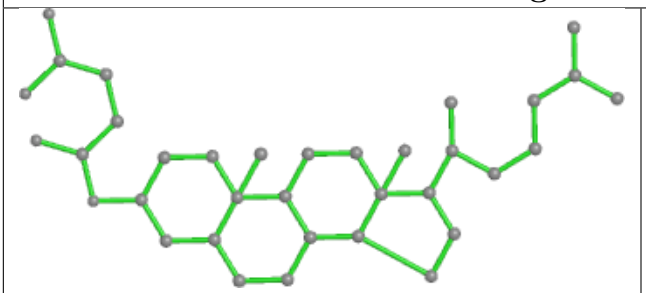
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	505	Y01	1	0
2	B	501	ACT	3	0
5	B	505	Y01	1	0
2	A	501	ACT	1	0
2	C	501	ACT	2	0
5	D	506	Y01	2	0
2	E	501	ACT	2	0

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 then 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.

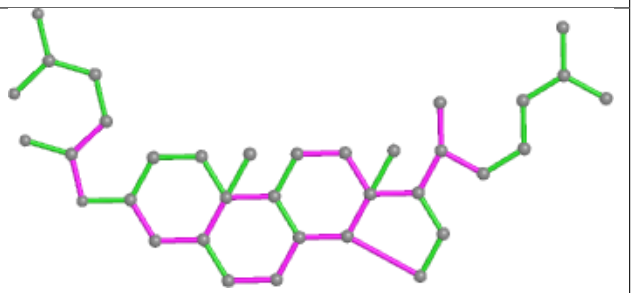
Ligand Y01 A 507



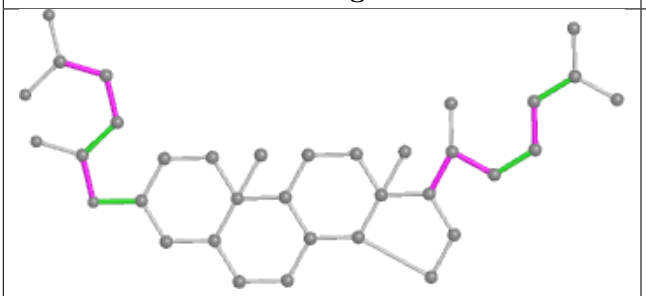
Ligand Y01 C 505



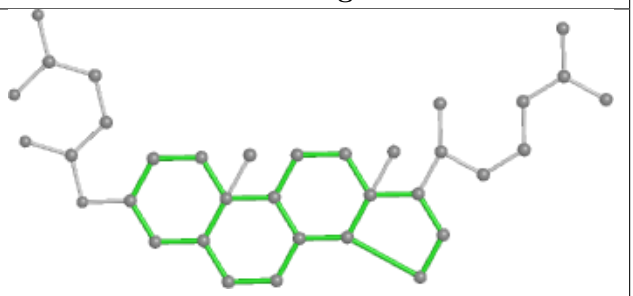
Bond lengths



Bond angles

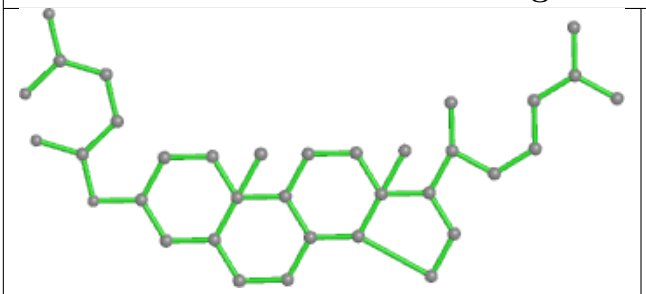


Torsions

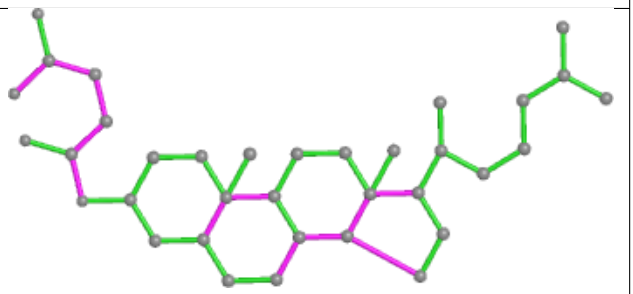


Rings

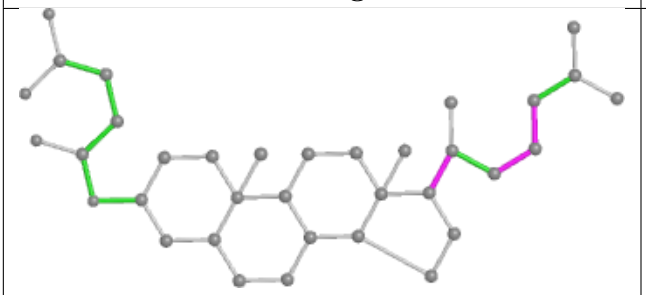
Ligand Y01 E 505



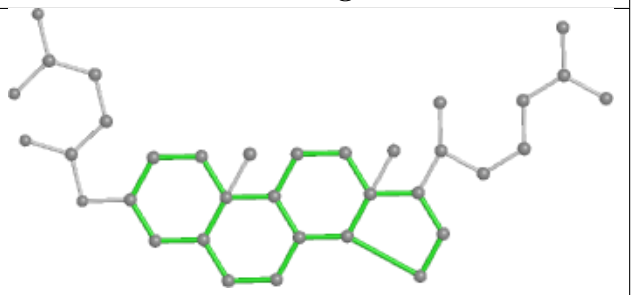
Bond lengths



Bond angles

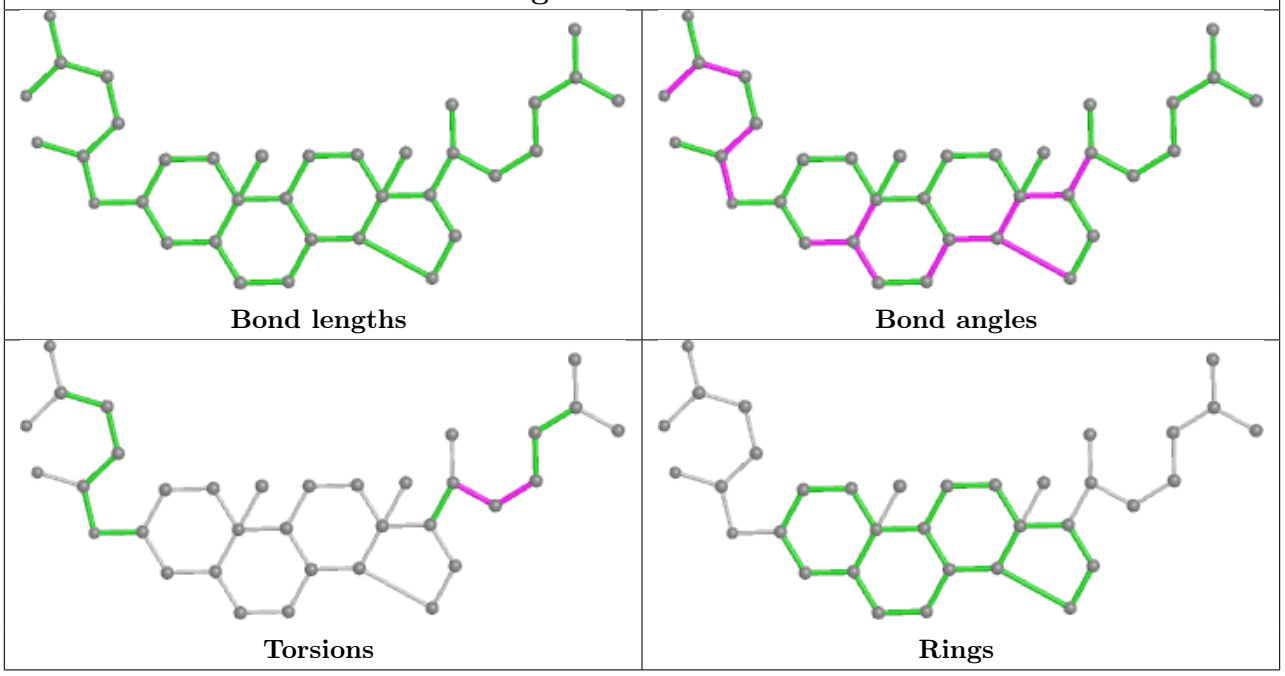


Torsions

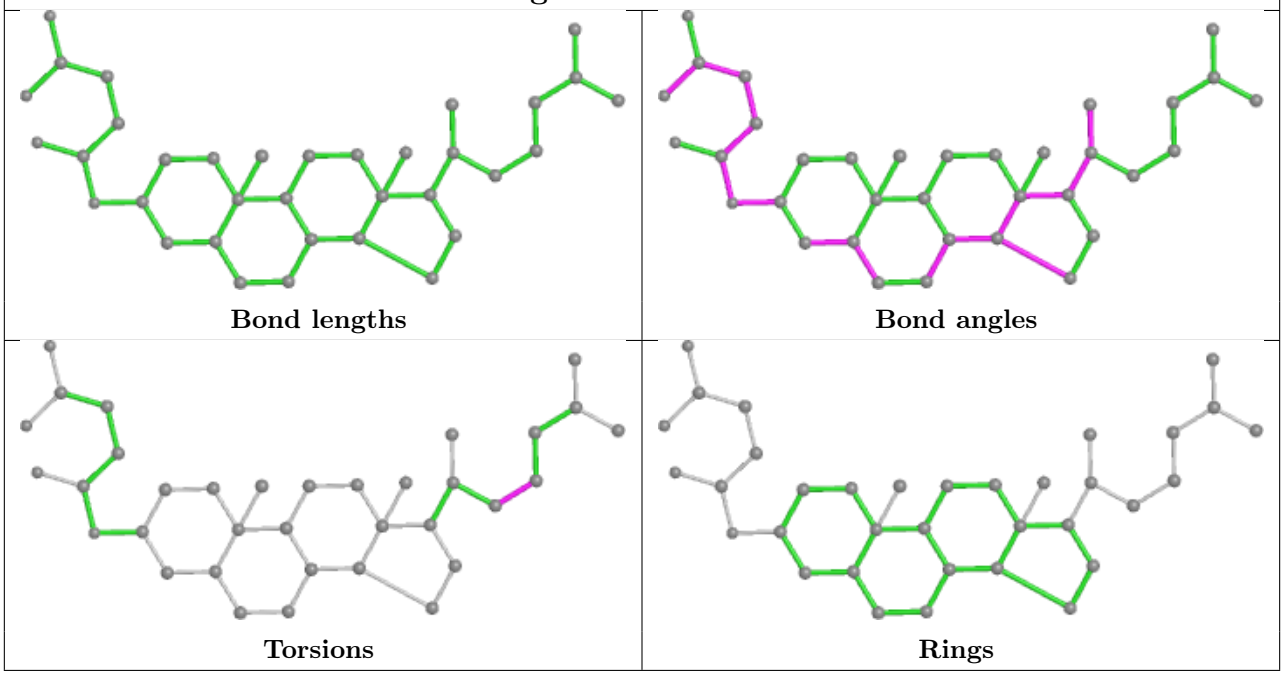


Rings

Ligand Y01 B 505



Ligand Y01 D 506



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	313/336 (93%)	0.48	12 (3%) 40 48	59, 84, 131, 217	0
1	B	313/336 (93%)	0.47	9 (2%) 51 61	58, 81, 121, 214	0
1	C	313/336 (93%)	0.36	7 (2%) 62 70	59, 84, 133, 262	0
1	D	313/336 (93%)	0.46	15 (4%) 30 36	59, 82, 116, 202	0
1	E	313/336 (93%)	0.40	7 (2%) 62 70	57, 80, 121, 218	0
All	All	1565/1680 (93%)	0.43	50 (3%) 47 56	57, 82, 124, 262	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	10	ILE	5.7
1	C	157	THR	4.8
1	A	12	ASP	4.1
1	B	32	LYS	4.0
1	E	60	VAL	4.0
1	C	60	VAL	3.8
1	D	11	ALA	3.7
1	A	11	ALA	3.5
1	B	147	LYS	3.4
1	A	135	ASP	3.1
1	D	60	VAL	3.1
1	A	157	THR	3.0
1	B	314	PRO	3.0
1	A	10	ILE	3.0
1	E	157	THR	3.0
1	B	113	LEU	3.0
1	D	157	THR	2.9
1	D	314	PRO	2.7
1	A	9	PRO	2.7
1	B	278	LYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	280	ALA	2.6
1	D	189	ILE	2.5
1	A	31	ASP	2.5
1	B	276	LEU	2.5
1	D	114	ASP	2.4
1	D	9	PRO	2.4
1	D	412	ALA	2.3
1	E	394	LEU	2.3
1	C	160	ASP	2.3
1	E	188	ARG	2.3
1	A	113	LEU	2.3
1	B	31	ASP	2.3
1	E	397	ILE	2.2
1	E	280	ALA	2.2
1	A	158	GLY	2.2
1	D	12	ASP	2.2
1	D	162	GLU	2.2
1	E	68	GLU	2.2
1	C	62	VAL	2.2
1	D	136	THR	2.2
1	C	32	LYS	2.2
1	A	32	LYS	2.1
1	A	115	PHE	2.1
1	B	65	TYR	2.1
1	B	279	VAL	2.1
1	C	193	TYR	2.0
1	C	278	LYS	2.0
1	A	134	VAL	2.0
1	D	193	TYR	2.0
1	D	188	ARG	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

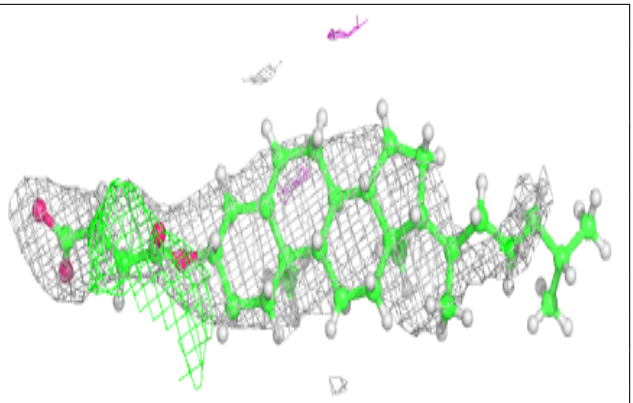
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

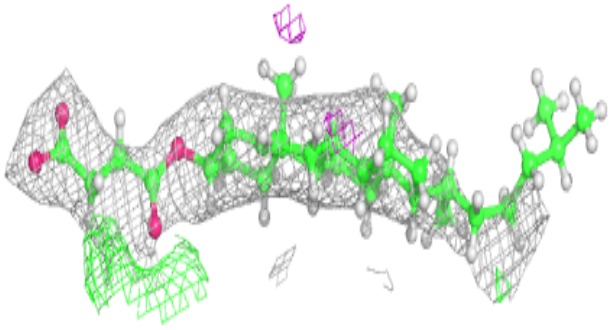
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	CL	A	505	1/1	0.70	0.29	116,116,116,116	0
4	D12	A	506	12/12	0.77	0.16	103,128,133,133	0
4	D12	B	503	12/12	0.78	0.40	111,137,149,149	0
4	D12	E	503	12/12	0.80	0.21	83,108,121,123	0
4	D12	E	504	12/12	0.80	0.52	105,131,150,150	0
5	Y01	D	506	35/35	0.80	0.44	116,166,200,200	0
3	CL	A	504	1/1	0.81	0.28	110,110,110,110	0
5	Y01	B	505	35/35	0.82	0.51	115,170,189,191	0
4	D12	D	504	12/12	0.82	0.25	100,124,132,133	0
4	D12	D	505	12/12	0.84	0.49	118,142,156,156	0
2	ACT	C	502	4/4	0.85	0.25	109,112,113,113	0
5	Y01	C	505	35/35	0.85	0.56	109,163,180,182	0
4	D12	B	504	12/12	0.85	0.17	85,114,125,126	0
5	Y01	E	505	35/35	0.86	0.30	116,150,168,170	0
5	Y01	A	507	35/35	0.88	0.42	125,161,174,175	0
4	D12	C	504	12/12	0.89	0.20	96,119,138,138	0
2	ACT	D	502	4/4	0.89	0.36	123,126,150,150	0
2	ACT	A	502	4/4	0.91	0.30	116,118,139,139	0
3	CL	D	503	1/1	0.94	0.15	71,71,71,71	0
3	CL	A	503	1/1	0.95	0.17	67,67,67,67	0
3	CL	B	502	1/1	0.95	0.16	76,76,76,76	0
2	ACT	A	501	4/4	0.97	0.18	67,77,77,80	0
3	CL	E	502	1/1	0.97	0.15	69,69,69,69	0
2	ACT	D	501	4/4	0.97	0.23	69,71,72,73	0
2	ACT	B	501	4/4	0.97	0.23	78,78,82,90	0
2	ACT	E	501	4/4	0.97	0.23	64,72,72,77	0
3	CL	C	503	1/1	0.98	0.16	73,73,73,73	0
2	ACT	C	501	4/4	0.98	0.28	74,74,78,84	0

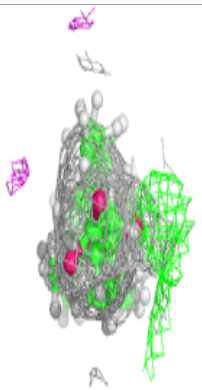
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around Y01 D 506:

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

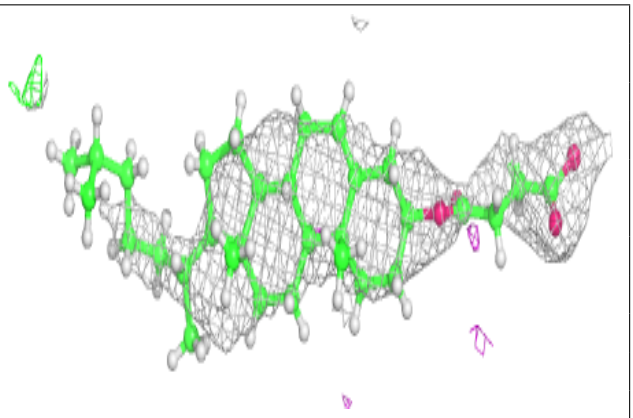


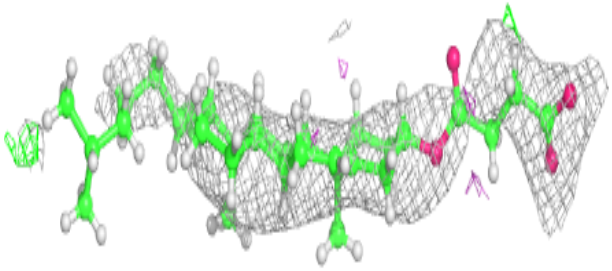


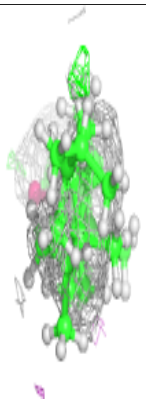


Electron density around Y01 B 505:

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

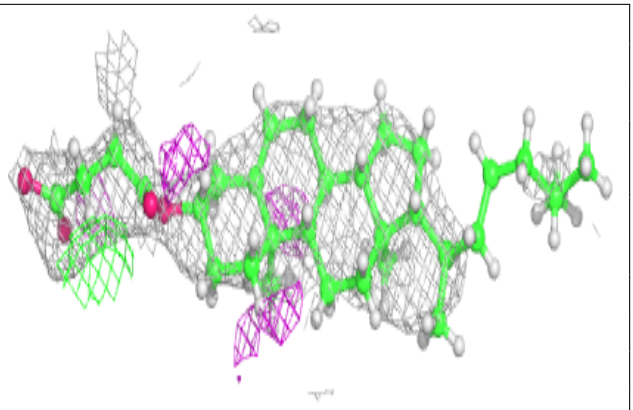


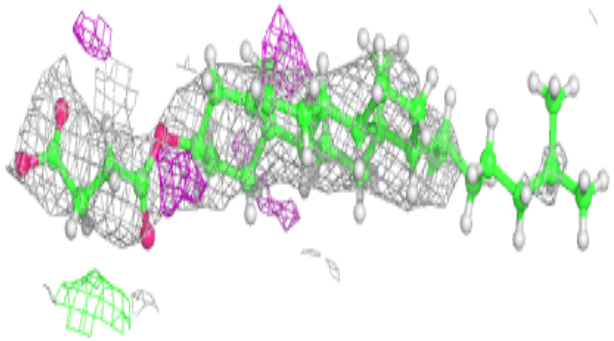


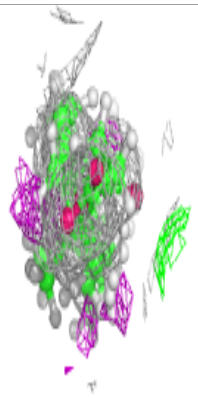


Electron density around Y01 C 505:

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

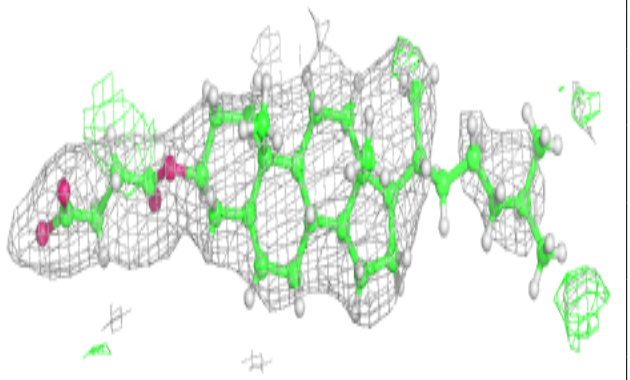


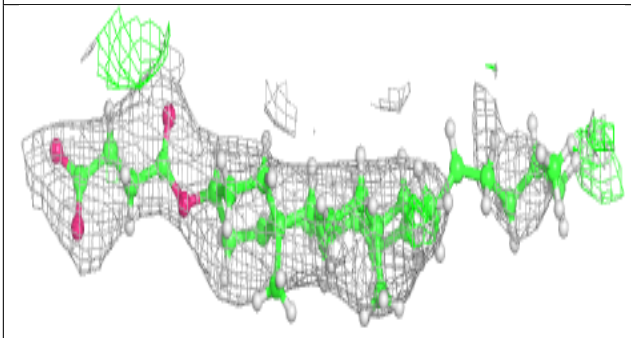


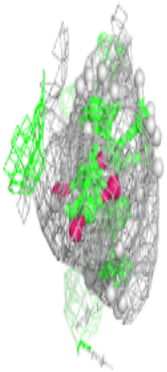


Electron density around Y01 E 505:

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

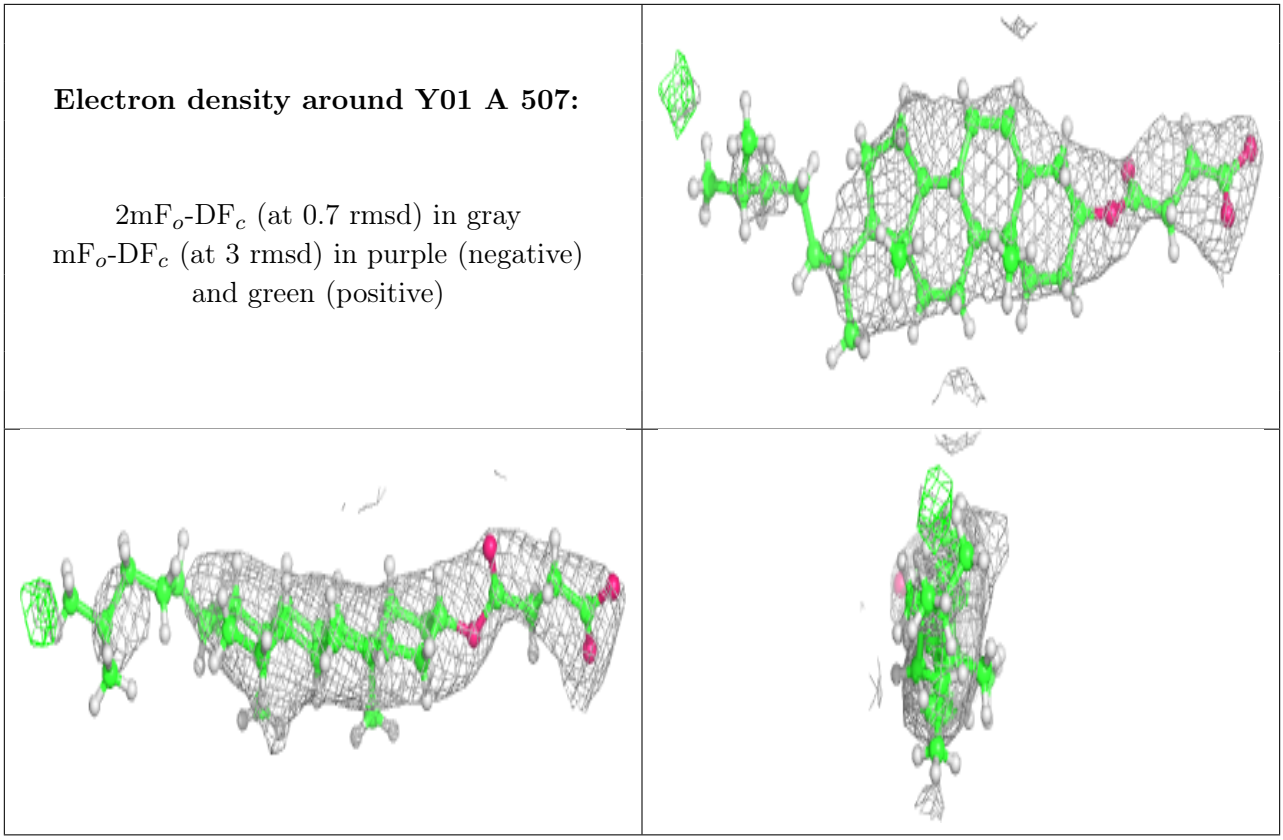






Electron density around Y01 A 507:

$2mF_o$-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.