



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:05 AM UTC

PDB ID : 1GSF / pdb_00001gsf
Title : GLUTATHIONE TRANSFERASE A1-1 COMPLEXED WITH
ETHACRYNIC ACID
Authors : L'Hermite, G.; Sinning, I.; Cameron, A.D.; Jones, T.A.
Deposited on : 1995-06-09
Resolution : 2.70 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	:	9.0.010 (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

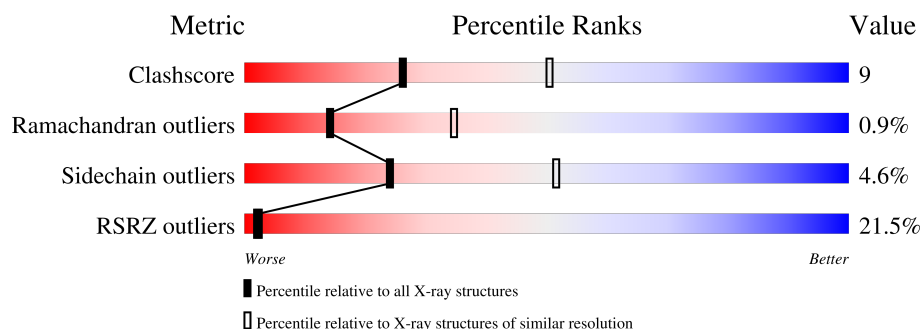
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	221	22% 77% 20% .
1	B	221	15% 77% 20% .
1	C	221	17% 77% 19% .
1	D	221	33% 76% 20% .

2 Entry composition [i](#)

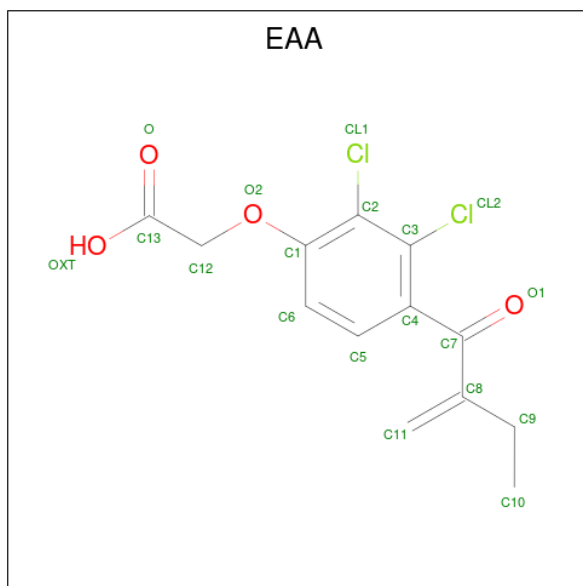
There are 3 unique types of molecules in this entry. The entry contains 7454 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GLUTATHIONE TRANSFERASE A1-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	1	0
			1807	1170	304	325	8			
1	B	221	Total	C	N	O	S	0	1	0
			1807	1170	304	325	8			
1	C	221	Total	C	N	O	S	0	1	0
			1807	1170	304	325	8			
1	D	221	Total	C	N	O	S	0	1	0
			1807	1170	304	325	8			

- Molecule 2 is ETHACRYNIC ACID (CCD ID: EAA) (formula: $C_{13}H_{12}Cl_2O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Cl	O	0	0
			19	13	2	4		
2	B	1	Total	C	Cl	O	0	0
			19	13	2	4		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	Cl	O	0	0
			19	13	2	4		
2	D	1	Total	C	Cl	O	0	0
			19	13	2	4		

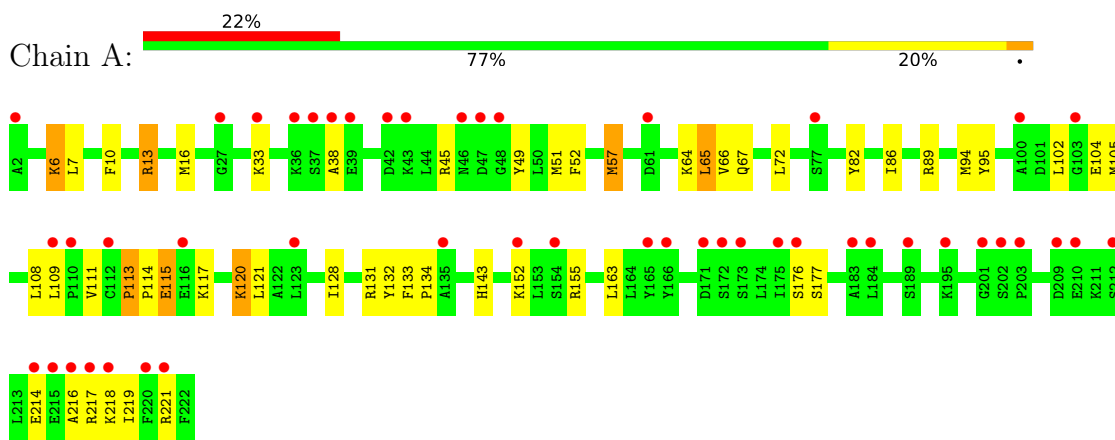
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	74	Total	O	0	0
			74	74		
3	B	76	Total	O	0	0
			76	76		

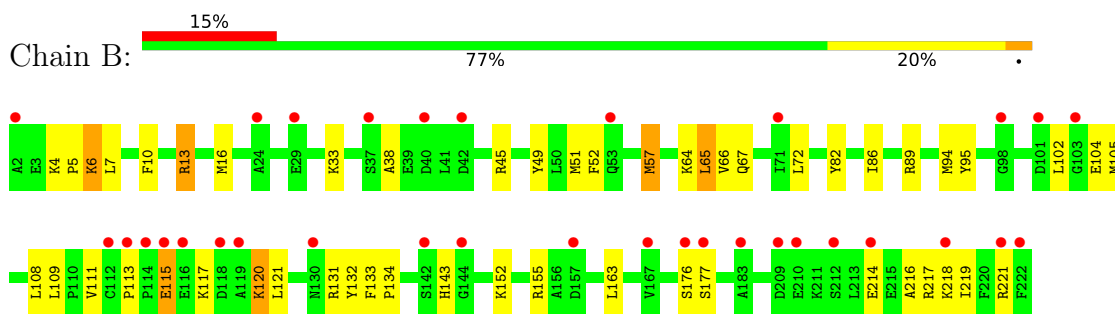
3 Residue-property plots

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.

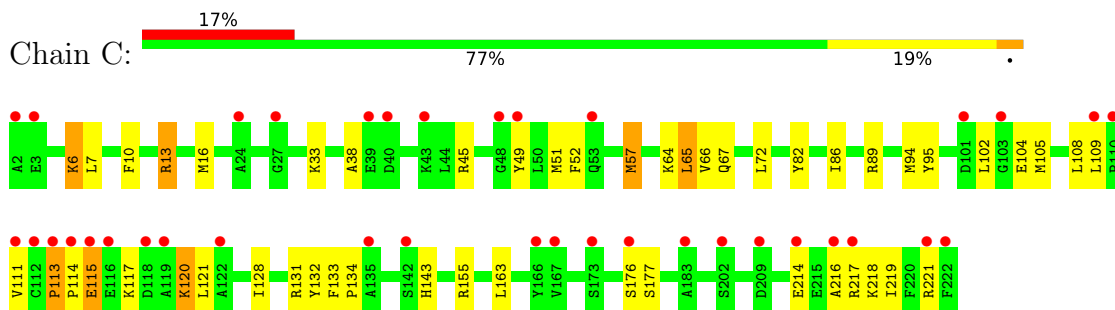
• Molecule 1: GLUTATHIONE TRANSFERASE A1-1



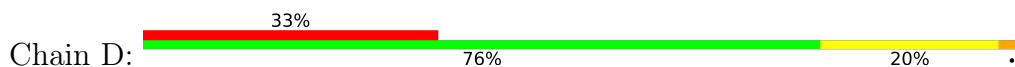
• Molecule 1: GLUTATHIONE TRANSFERASE A1-1

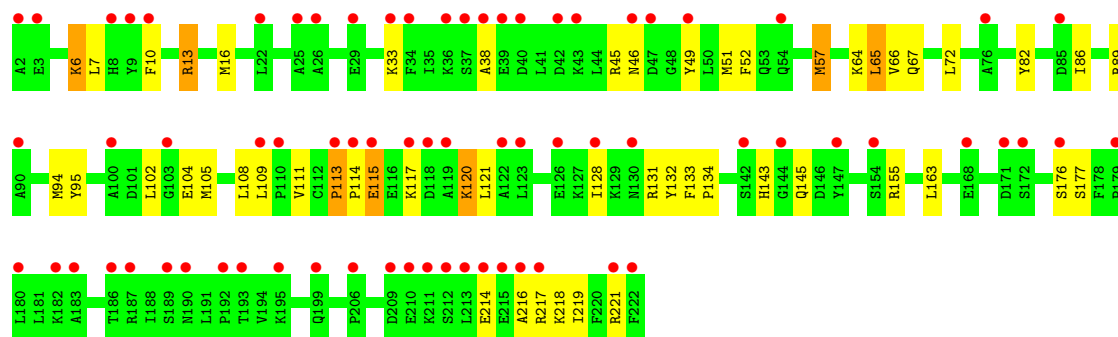


• Molecule 1: GLUTATHIONE TRANSFERASE A1-1



• Molecule 1: GLUTATHIONE TRANSFERASE A1-1





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	101.80Å 95.60Å 105.30Å 90.00° 92.20° 90.00°	Depositor
Resolution (Å)	7.50 – 2.70 7.50 – 2.70	Depositor EDS
% Data completeness (in resolution range)	94.9 (7.50-2.70) 89.6 (7.50-2.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.99 (at 2.71Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.229 , 0.261 0.312 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	44.0	Xtriage
Anisotropy	0.356	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 53.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.012 for -h,-k,l	Xtriage
F_o, F_c correlation	0.81	EDS
Total number of atoms	7454	wwPDB-VP
Average B, all atoms (Å ²)	43.0	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 34.76 % 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 6.4903e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EAA

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	0.64	0/1843	1.02	6/2475 (0.2%)
1	B	0.64	0/1843	1.02	6/2475 (0.2%)
1	C	0.64	0/1843	1.02	6/2475 (0.2%)
1	D	0.64	0/1843	1.02	6/2475 (0.2%)
All	All	0.64	0/7372	1.02	24/9900 (0.2%)

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	113	PRO	CA-C-N	6.95	128.53	119.84
1	C	113	PRO	C-N-CA	6.95	128.53	119.84
1	A	113	PRO	CA-C-N	6.93	128.51	119.84
1	A	113	PRO	C-N-CA	6.93	128.51	119.84
1	B	113	PRO	CA-C-N	6.93	128.50	119.84
1	B	113	PRO	C-N-CA	6.93	128.50	119.84
1	D	113	PRO	CA-C-N	6.92	128.49	119.84
1	D	113	PRO	C-N-CA	6.92	128.49	119.84
1	C	131	ARG	N-CA-C	5.67	119.72	112.12
1	A	131	ARG	N-CA-C	5.63	119.67	112.12
1	B	131	ARG	N-CA-C	5.62	119.66	112.12
1	D	131	ARG	N-CA-C	5.61	119.64	112.12
1	A	132	TYR	N-CA-C	5.45	118.05	111.40
1	C	132	TYR	N-CA-C	5.42	118.02	111.40
1	D	132	TYR	N-CA-C	5.42	118.01	111.40
1	B	132	TYR	N-CA-C	5.42	118.01	111.40
1	A	57	MET	N-CA-C	-5.32	100.98	109.23
1	C	57	MET	N-CA-C	-5.31	100.99	109.23
1	D	57	MET	N-CA-C	-5.29	101.03	109.23
1	B	57	MET	N-CA-C	-5.29	101.04	109.23

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	C	176	SER	N-CA-C	5.21	117.64	111.33
1	A	176	SER	N-CA-C	5.21	117.64	111.33
1	D	176	SER	N-CA-C	5.20	117.62	111.33
1	B	176	SER	N-CA-C	5.18	117.60	111.33

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	1807	0	1875	46	0
1	B	1807	0	1875	45	0
1	C	1807	0	1875	45	0
1	D	1807	0	1875	44	1
2	A	19	0	11	0	0
2	B	19	0	11	0	0
2	C	19	0	11	0	0
2	D	19	0	11	0	0
3	A	74	0	0	1	0
3	B	76	0	0	1	0
All	All	7454	0	7544	139	1

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 9.

All (139) 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:89[B]:ARG:CZ	1:B:89[B]:ARG:HH22	1.07	1.65
1:C:89[B]:ARG:HH22	1:D:89[B]:ARG:CZ	1.05	1.59
1:C:89[B]:ARG:CZ	1:D:89[B]:ARG:HH22	1.16	1.57
1:A:89[B]:ARG:HH22	1:B:89[B]:ARG:CZ	1.19	1.51
1:C:89[B]:ARG:HH12	1:D:89[B]:ARG:NH1	1.29	1.31
1:C:89[B]:ARG:NH1	1:D:89[B]:ARG:HH12	1.26	1.29

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89[B]:ARG:HH12	1:B:89[B]:ARG:NH1	1.28	1.29
1:A:89[B]:ARG:NH1	1:B:89[B]:ARG:HH12	1.32	1.26
1:C:89[B]:ARG:CZ	1:D:89[B]:ARG:NH2	1.96	0.88
1:C:89[B]:ARG:NH1	1:D:89[B]:ARG:NH1	2.02	0.85
1:A:89[B]:ARG:HH21	1:B:89[B]:ARG:NH2	1.33	0.82
1:A:89[B]:ARG:NH1	1:B:89[B]:ARG:NH1	2.05	0.81
1:C:89[B]:ARG:NH2	1:D:89[B]:ARG:CZ	1.87	0.80
1:A:89[B]:ARG:NH2	1:B:89[B]:ARG:CZ	1.99	0.80
1:C:89[B]:ARG:NH2	1:D:89[B]:ARG:HH21	1.31	0.79
1:A:102:LEU:HA	1:A:105:MET:HE3	1.66	0.78
1:B:102:LEU:HA	1:B:105:MET:HE3	1.66	0.78
1:C:102:LEU:HA	1:C:105:MET:HE3	1.66	0.77
1:D:102:LEU:HA	1:D:105:MET:HE3	1.66	0.76
1:A:89[B]:ARG:CZ	1:B:89[B]:ARG:NH2	1.90	0.73
1:A:89[B]:ARG:NH2	1:B:89[B]:ARG:HH21	1.21	0.71
1:B:45:ARG:HH12	1:B:221:ARG:HH22	1.41	0.69
1:C:89[B]:ARG:NH2	1:D:89[B]:ARG:NH2	0.73	0.68
1:A:45:ARG:HH12	1:A:221:ARG:HH22	1.41	0.67
1:A:89[B]:ARG:NH2	1:B:89[B]:ARG:HH22	0.35	0.67
1:D:45:ARG:HH12	1:D:221:ARG:HH22	1.41	0.67
1:C:89[B]:ARG:HH21	1:D:89[B]:ARG:NH2	1.19	0.67
1:C:45:ARG:HH12	1:C:221:ARG:HH22	1.41	0.67
1:A:89[B]:ARG:NH2	1:B:89[B]:ARG:NH2	0.76	0.66
1:C:89[B]:ARG:HH22	1:D:89[B]:ARG:NH2	0.35	0.65
1:C:89[B]:ARG:CZ	1:D:89[B]:ARG:HH12	2.07	0.64
1:C:89[B]:ARG:CZ	1:D:89[B]:ARG:NH1	2.60	0.63
1:C:109:LEU:HD21	1:C:121:LEU:HD13	1.80	0.63
1:A:109:LEU:HD21	1:A:121:LEU:HD13	1.80	0.63
1:D:109:LEU:HD21	1:D:121:LEU:HD13	1.80	0.63
1:B:109:LEU:HD21	1:B:121:LEU:HD13	1.80	0.62
1:C:89[B]:ARG:HH12	1:D:89[B]:ARG:HH12	0.66	0.60
1:A:95:TYR:HE1	1:B:51:MET:HE1	1.66	0.59
1:A:89[B]:ARG:HH22	1:B:89[B]:ARG:NH2	0.19	0.59
1:A:89[B]:ARG:HH12	1:B:89[B]:ARG:HH12	0.67	0.59
1:A:89[B]:ARG:NH1	1:B:89[B]:ARG:CZ	2.63	0.59
1:C:89[B]:ARG:CZ	1:D:89[B]:ARG:CZ	2.63	0.57
1:C:89[B]:ARG:NH2	1:D:89[B]:ARG:HH22	0.20	0.55
1:A:89[B]:ARG:HH22	1:B:89[B]:ARG:HH21	0.98	0.55
1:D:214:GLU:OE2	1:D:217:ARG:HD3	2.07	0.55
1:A:51:MET:HE1	1:B:95:TYR:HE1	1.70	0.55
1:B:214:GLU:OE2	1:B:217:ARG:HD3	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:GLU:OE2	1:A:217:ARG:HD3	2.07	0.55
1:B:82:TYR:CD1	1:B:89[A]:ARG:HG2	2.42	0.55
1:A:82:TYR:CD1	1:A:89[A]:ARG:HG2	2.42	0.54
1:C:82:TYR:CD1	1:C:89[A]:ARG:HG2	2.42	0.54
1:C:214:GLU:OE2	1:C:217:ARG:HD3	2.07	0.54
1:D:82:TYR:CD1	1:D:89[A]:ARG:HG2	2.42	0.54
1:C:94:MET:HE2	1:D:65:LEU:HD12	1.90	0.54
1:C:217:ARG:O	1:C:221:ARG:HA	2.08	0.54
1:B:115:GLU:H	1:B:115:GLU:CD	2.16	0.54
1:A:65:LEU:HD12	1:B:94:MET:HE2	1.90	0.53
1:A:89[B]:ARG:HH12	1:B:89[B]:ARG:CZ	2.09	0.53
1:A:115:GLU:H	1:A:115:GLU:CD	2.16	0.53
1:A:217:ARG:O	1:A:221:ARG:HA	2.08	0.53
1:D:217:ARG:O	1:D:221:ARG:HA	2.08	0.53
1:B:217:ARG:O	1:B:221:ARG:HA	2.08	0.53
1:C:95:TYR:HE1	1:D:51:MET:HE1	1.73	0.53
1:C:111:VAL:O	1:C:111:VAL:HG12	2.09	0.53
1:C:115:GLU:H	1:C:115:GLU:CD	2.16	0.53
1:D:115:GLU:CD	1:D:115:GLU:H	2.17	0.53
1:D:111:VAL:HG12	1:D:111:VAL:O	2.09	0.52
1:B:111:VAL:O	1:B:111:VAL:HG12	2.09	0.52
1:C:51:MET:HE1	1:D:95:TYR:HE1	1.74	0.52
1:B:66:VAL:O	1:B:67:GLN:HB2	2.11	0.51
1:A:66:VAL:O	1:A:67:GLN:HB2	2.11	0.50
1:A:111:VAL:HG12	1:A:111:VAL:O	2.09	0.50
1:B:152:LYS:HE2	3:B:284:HOH:O	2.11	0.50
1:A:152:LYS:HE2	3:A:279:HOH:O	2.11	0.50
1:D:49:TYR:HB3	1:D:57:MET:HE2	1.94	0.50
1:D:66:VAL:O	1:D:67:GLN:HB2	2.11	0.49
1:B:7:LEU:HB3	1:B:16:MET:HE3	1.94	0.49
1:C:66:VAL:O	1:C:67:GLN:HB2	2.11	0.49
1:B:49:TYR:HB3	1:B:57:MET:HE2	1.94	0.49
1:C:7:LEU:HB3	1:C:16:MET:HE3	1.94	0.49
1:D:7:LEU:HB3	1:D:16:MET:HE3	1.94	0.49
1:A:7:LEU:HB3	1:A:16:MET:HE3	1.94	0.49
1:C:49:TYR:HB3	1:C:57:MET:HE2	1.94	0.48
1:A:49:TYR:HB3	1:A:57:MET:HE2	1.94	0.48
1:C:113:PRO:HA	1:C:114:PRO:HD2	1.69	0.47
1:C:82:TYR:CE1	1:C:89[B]:ARG:HG2	2.50	0.47
1:D:10:PHE:CE2	1:D:216:ALA:HB2	2.50	0.47
1:B:82:TYR:CE1	1:B:89[B]:ARG:HG2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:PHE:CE2	1:A:216:ALA:HB2	2.50	0.46
1:B:10:PHE:CE2	1:B:216:ALA:HB2	2.50	0.46
1:A:82:TYR:CE1	1:A:89[B]:ARG:HG2	2.50	0.46
1:A:94:MET:HE2	1:B:65:LEU:HD12	1.98	0.46
1:D:82:TYR:CE1	1:D:89[B]:ARG:HG2	2.50	0.46
1:C:10:PHE:CE2	1:C:216:ALA:HB2	2.50	0.45
1:A:102:LEU:HD23	1:A:163:LEU:HD21	1.99	0.44
1:D:38:ALA:HA	1:D:219:ILE:HG23	1.99	0.44
1:C:102:LEU:HD23	1:C:163:LEU:HD21	1.99	0.44
1:A:38:ALA:HA	1:A:219:ILE:HG23	1.99	0.44
1:A:89[B]:ARG:HH21	1:B:89[B]:ARG:HH21	1.18	0.44
1:C:104:GLU:O	1:C:108:LEU:HG	2.18	0.44
1:C:38:ALA:HA	1:C:219:ILE:HG23	1.99	0.44
1:D:102:LEU:HD23	1:D:163:LEU:HD21	1.99	0.44
1:B:38:ALA:HA	1:B:219:ILE:HG23	1.99	0.44
1:B:102:LEU:HD23	1:B:163:LEU:HD21	1.99	0.43
1:A:104:GLU:O	1:A:108:LEU:HG	2.18	0.43
1:A:113:PRO:HA	1:A:114:PRO:HD2	1.69	0.43
1:B:104:GLU:O	1:B:108:LEU:HG	2.18	0.43
1:B:120:LYS:NZ	1:B:120:LYS:HB3	2.33	0.43
1:D:104:GLU:O	1:D:108:LEU:HG	2.18	0.43
1:C:52:PHE:HD2	1:C:67:GLN:HE21	1.66	0.43
1:B:52:PHE:HD2	1:B:67:GLN:HE21	1.67	0.43
1:D:113:PRO:HA	1:D:114:PRO:HD2	1.69	0.43
1:B:72:LEU:HB3	1:B:155:ARG:NH2	2.34	0.43
1:D:72:LEU:HB3	1:D:155:ARG:NH2	2.34	0.43
1:C:72:LEU:HB3	1:C:155:ARG:NH2	2.34	0.42
1:A:120:LYS:HB3	1:A:120:LYS:NZ	2.34	0.42
1:D:120:LYS:HB3	1:D:120:LYS:NZ	2.33	0.42
1:C:6:LYS:HG2	1:C:33:LYS:HE3	2.02	0.42
1:D:52:PHE:HD2	1:D:67:GLN:HE21	1.67	0.42
1:C:120:LYS:NZ	1:C:120:LYS:HB3	2.34	0.42
1:D:6:LYS:HG2	1:D:33:LYS:HE3	2.02	0.41
1:A:52:PHE:HD2	1:A:67:GLN:HE21	1.67	0.41
1:A:72:LEU:HB3	1:A:155:ARG:NH2	2.34	0.41
1:D:49:TYR:CE1	1:D:64:LYS:HE2	2.56	0.41
1:A:6:LYS:HG2	1:A:33:LYS:HE3	2.02	0.41
1:B:133:PHE:N	1:B:134:PRO:CD	2.84	0.41
1:C:65:LEU:HD12	1:D:94:MET:HE2	2.03	0.41
1:C:102:LEU:HD21	1:C:128:ILE:HG12	2.03	0.41
1:A:133:PHE:N	1:A:134:PRO:CD	2.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:6:LYS:HG2	1:B:33:LYS:HE3	2.02	0.41
1:B:49:TYR:CE1	1:B:64:LYS:HE2	2.56	0.41
1:D:102:LEU:HD21	1:D:128:ILE:HG12	2.03	0.41
1:A:102:LEU:HD21	1:A:128:ILE:HG12	2.03	0.40
1:C:133:PHE:N	1:C:134:PRO:CD	2.84	0.40
1:A:49:TYR:CE1	1:A:64:LYS:HE2	2.56	0.40
1:D:133:PHE:N	1:D:134:PRO:CD	2.84	0.40
1:C:38:ALA:CA	1:C:219:ILE:HG23	2.52	0.40
1:C:49:TYR:CE1	1:C:64:LYS:HE2	2.56	0.40
1:B:4:LYS:HA	1:B:5:PRO:HD3	1.89	0.40

All (1) 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:D:46:ASN:O	1:D:145:GLN:OE1[4_656]	1.66	0.54

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	220/221 (100%)	209 (95%)	9 (4%)	2 (1%)	14	35
1	B	220/221 (100%)	209 (95%)	9 (4%)	2 (1%)	14	35
1	C	220/221 (100%)	209 (95%)	9 (4%)	2 (1%)	14	35
1	D	220/221 (100%)	209 (95%)	9 (4%)	2 (1%)	14	35
All	All	880/884 (100%)	836 (95%)	36 (4%)	8 (1%)	14	35

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	117	LYS
1	B	117	LYS
1	C	117	LYS
1	D	117	LYS
1	A	13	ARG
1	B	13	ARG
1	C	13	ARG
1	D	13	ARG

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	196/195 (100%)	187 (95%)	9 (5%)	24	51	
1	B	196/195 (100%)	187 (95%)	9 (5%)	24	51	
1	C	196/195 (100%)	187 (95%)	9 (5%)	24	51	
1	D	196/195 (100%)	187 (95%)	9 (5%)	24	51	
All	All	784/780 (100%)	748 (95%)	36 (5%)	24	51	

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LYS
1	A	13	ARG
1	A	65	LEU
1	A	86	ILE
1	A	115	GLU
1	A	120	LYS
1	A	143	HIS
1	A	177	SER
1	A	218	LYS
1	B	6	LYS
1	B	13	ARG
1	B	65	LEU
1	B	86	ILE
1	B	115	GLU

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Mol	Chain	Res	Type
1	B	120	LYS
1	B	143	HIS
1	B	177	SER
1	B	218	LYS
1	C	6	LYS
1	C	13	ARG
1	C	65	LEU
1	C	86	ILE
1	C	115	GLU
1	C	120	LYS
1	C	143	HIS
1	C	177	SER
1	C	218	LYS
1	D	6	LYS
1	D	13	ARG
1	D	65	LEU
1	D	86	ILE
1	D	115	GLU
1	D	120	LYS
1	D	143	HIS
1	D	177	SER
1	D	218	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	GLN
1	A	67	GLN
1	A	130	ASN
1	A	199	GLN
1	B	53	GLN
1	B	67	GLN
1	B	130	ASN
1	B	199	GLN
1	C	53	GLN
1	C	67	GLN
1	C	130	ASN
1	C	199	GLN
1	D	53	GLN
1	D	67	GLN
1	D	130	ASN
1	D	199	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 ⓘ

4 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$
2	EAA	A	223	-	18,19,19	1.75	2 (11%)	21,26,26	1.23	2 (9%)
2	EAA	B	223	-	18,19,19	1.76	2 (11%)	21,26,26	1.24	2 (9%)
2	EAA	C	223	-	18,19,19	1.75	2 (11%)	21,26,26	1.24	2 (9%)
2	EAA	D	223	-	18,19,19	1.75	2 (11%)	21,26,26	1.23	2 (9%)

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
2	EAA	A	223	-	-	5/15/15/15	0/1/1/1
2	EAA	B	223	-	-	5/15/15/15	0/1/1/1
2	EAA	C	223	-	-	5/15/15/15	0/1/1/1
2	EAA	D	223	-	-	5/15/15/15	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	223	EAA	C11-C8	-5.22	1.22	1.32
2	A	223	EAA	C11-C8	-5.21	1.22	1.32
2	C	223	EAA	C11-C8	-5.19	1.22	1.32
2	D	223	EAA	C11-C8	-5.18	1.22	1.32
2	C	223	EAA	C4-C7	-2.87	1.43	1.49
2	B	223	EAA	C4-C7	-2.86	1.43	1.49
2	A	223	EAA	C4-C7	-2.85	1.43	1.49
2	D	223	EAA	C4-C7	-2.85	1.43	1.49

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	223	EAA	O2-C1-C2	3.33	121.11	115.77
2	A	223	EAA	O2-C1-C2	3.31	121.09	115.77
2	D	223	EAA	O2-C1-C2	3.31	121.08	115.77
2	C	223	EAA	O2-C1-C2	3.31	121.08	115.77
2	C	223	EAA	C1-C2-C3	2.81	121.72	119.86
2	B	223	EAA	C1-C2-C3	2.78	121.70	119.86
2	A	223	EAA	C1-C2-C3	2.76	121.69	119.86
2	D	223	EAA	C1-C2-C3	2.74	121.67	119.86

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	223	EAA	C4-C7-C8-C9
2	A	223	EAA	C4-C7-C8-C11
2	A	223	EAA	O1-C7-C8-C9
2	A	223	EAA	O1-C7-C8-C11
2	A	223	EAA	C7-C8-C9-C10
2	B	223	EAA	C4-C7-C8-C9
2	B	223	EAA	C4-C7-C8-C11
2	B	223	EAA	O1-C7-C8-C9
2	B	223	EAA	O1-C7-C8-C11
2	B	223	EAA	C7-C8-C9-C10
2	C	223	EAA	C4-C7-C8-C9
2	C	223	EAA	C4-C7-C8-C11
2	C	223	EAA	O1-C7-C8-C9
2	C	223	EAA	O1-C7-C8-C11
2	C	223	EAA	C7-C8-C9-C10
2	D	223	EAA	C4-C7-C8-C9

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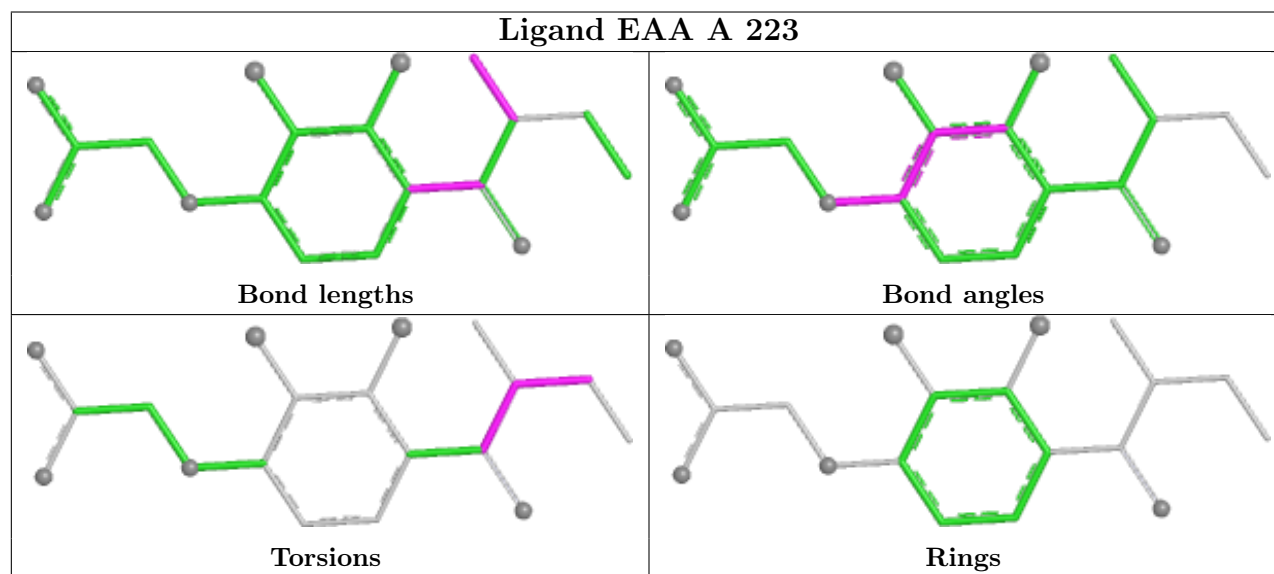
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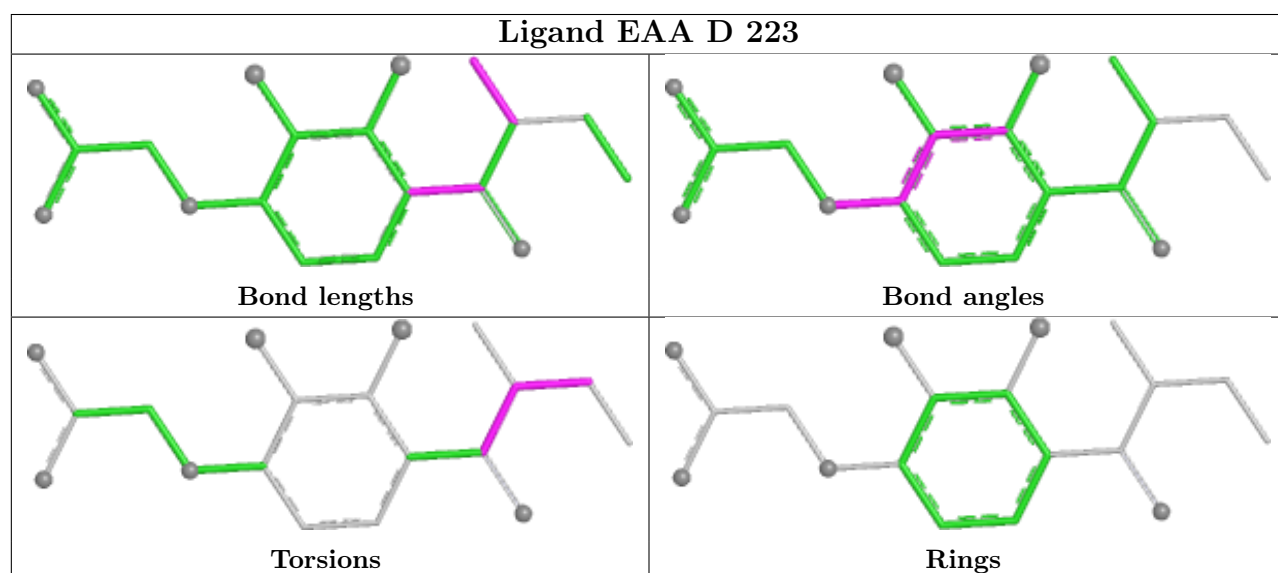
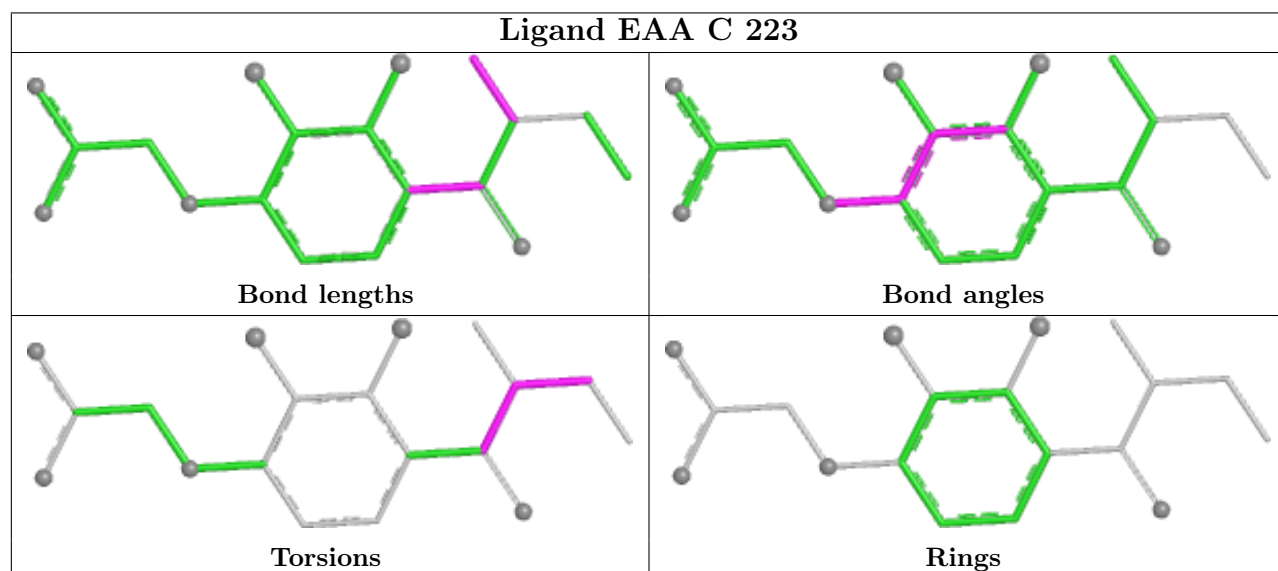
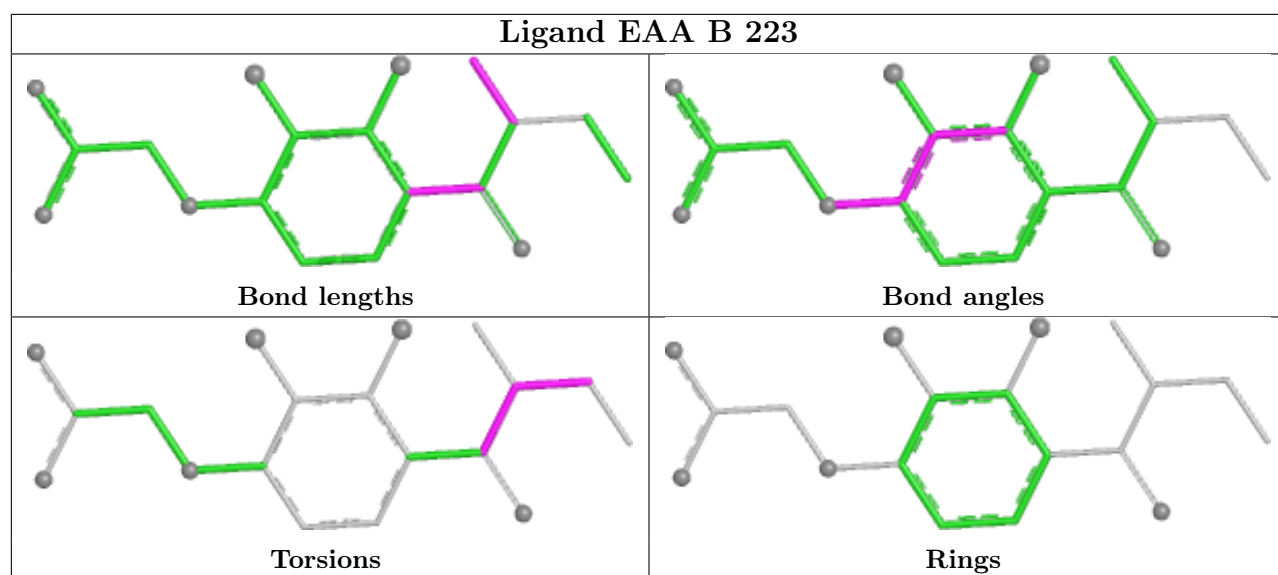
Mol	Chain	Res	Type	Atoms
2	D	223	EAA	C4-C7-C8-C11
2	D	223	EAA	O1-C7-C8-C9
2	D	223	EAA	O1-C7-C8-C11
2	D	223	EAA	C7-C8-C9-C10

There are no ring outliers.

No monomer is involved in short contacts.

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	221/221 (100%)	1.32	48 (21%) 2 2	7, 35, 85, 118	1 (0%)
1	B	221/221 (100%)	1.13	33 (14%) 5 4	7, 35, 85, 118	1 (0%)
1	C	221/221 (100%)	1.10	37 (16%) 4 3	7, 35, 85, 118	1 (0%)
1	D	221/221 (100%)	1.54	72 (32%) 1 1	7, 35, 85, 118	1 (0%)
All	All	884/884 (100%)	1.27	190 (21%) 2 2	7, 36, 88, 118	4 (0%)

All (190) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	118	ASP	5.4
1	A	42	ASP	5.1
1	C	115	GLU	5.0
1	D	209	ASP	4.9
1	D	40	ASP	4.9
1	A	209	ASP	4.6
1	B	119	ALA	4.5
1	C	118	ASP	4.5
1	D	38	ALA	4.4
1	D	42	ASP	4.3
1	D	213	LEU	4.3
1	D	211	LYS	4.3
1	B	115	GLU	4.3
1	C	103	GLY	4.2
1	C	217	ARG	4.1
1	D	33	LYS	4.1
1	C	119	ALA	4.1
1	D	114	PRO	4.0
1	A	214	GLU	3.9
1	C	214	GLU	3.9
1	B	113	PRO	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	39	GLU	3.8
1	D	189	SER	3.8
1	A	2	ALA	3.8
1	B	24	ALA	3.8
1	C	114	PRO	3.8
1	D	29	GLU	3.7
1	D	37	SER	3.7
1	D	47	ASP	3.6
1	D	214	GLU	3.6
1	A	210	GLU	3.6
1	D	168	GLU	3.6
1	C	221	ARG	3.5
1	C	110	PRO	3.5
1	B	176	SER	3.5
1	C	173	SER	3.4
1	A	176	SER	3.4
1	B	112	CYS	3.4
1	D	183	ALA	3.4
1	D	212	SER	3.4
1	B	144	GLY	3.4
1	D	206	PRO	3.4
1	C	222	PHE	3.4
1	B	212	SER	3.3
1	A	43	LYS	3.3
1	C	112	CYS	3.3
1	D	142	SER	3.3
1	D	217	ARG	3.2
1	D	118	ASP	3.2
1	A	33	LYS	3.2
1	A	183	ALA	3.2
1	A	217	ARG	3.2
1	A	195	LYS	3.1
1	B	116	GLU	3.1
1	B	221	ARG	3.1
1	D	154	SER	3.1
1	D	113	PRO	3.0
1	C	142	SER	3.0
1	A	215	GLU	3.0
1	A	216	ALA	3.0
1	D	119	ALA	3.0
1	D	179	PRO	2.9
1	B	209	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	46	ASN	2.9
1	D	171	ASP	2.9
1	A	38	ALA	2.9
1	A	116	GLU	2.9
1	A	202	SER	2.9
1	D	187	ARG	2.9
1	D	126	GLU	2.9
1	D	221	ARG	2.9
1	A	37	SER	2.9
1	B	103	GLY	2.9
1	D	123	LEU	2.9
1	D	195	LYS	2.9
1	D	215	GLU	2.8
1	D	192	PRO	2.8
1	C	202	SER	2.8
1	D	46	ASN	2.8
1	D	182	LYS	2.8
1	B	40	ASP	2.8
1	D	190	ASN	2.8
1	A	220	PHE	2.8
1	B	214	GLU	2.8
1	D	222	PHE	2.8
1	A	212	SER	2.8
1	B	130	ASN	2.8
1	A	171	ASP	2.8
1	A	39	GLU	2.8
1	A	110	PRO	2.7
1	B	218	LYS	2.7
1	D	109	LEU	2.7
1	C	176	SER	2.7
1	D	54	GLN	2.7
1	A	218	LYS	2.7
1	D	103	GLY	2.7
1	A	100	ALA	2.7
1	B	157	ASP	2.7
1	C	113	PRO	2.7
1	C	167	VAL	2.6
1	C	40	ASP	2.6
1	C	27	GLY	2.6
1	A	61	ASP	2.6
1	C	48	GLY	2.6
1	C	109	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	48	GLY	2.5
1	C	135	ALA	2.5
1	B	210	GLU	2.5
1	D	110	PRO	2.5
1	C	2	ALA	2.5
1	C	24	ALA	2.5
1	A	154	SER	2.5
1	C	209	ASP	2.5
1	D	43	LYS	2.4
1	A	173	SER	2.4
1	B	42	ASP	2.4
1	D	117	LYS	2.4
1	A	184	LEU	2.4
1	D	172	SER	2.4
1	B	53	GLN	2.4
1	A	166	TYR	2.4
1	C	39	GLU	2.4
1	D	3	GLU	2.4
1	B	114	PRO	2.4
1	A	109	LEU	2.3
1	C	116	GLU	2.3
1	D	210	GLU	2.3
1	D	10	PHE	2.3
1	D	22	LEU	2.3
1	A	36	LYS	2.3
1	B	2	ALA	2.3
1	D	9	TYR	2.3
1	B	101	ASP	2.3
1	A	103	GLY	2.3
1	C	53	GLN	2.3
1	D	130	ASN	2.3
1	A	135	ALA	2.2
1	D	128	ILE	2.2
1	A	189	SER	2.2
1	B	37	SER	2.2
1	B	222	PHE	2.2
1	D	2	ALA	2.2
1	D	8	HIS	2.2
1	D	36	LYS	2.2
1	A	112	CYS	2.2
1	D	186	THR	2.2
1	A	203	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	142	SER	2.2
1	D	176	SER	2.2
1	B	183	ALA	2.2
1	D	85	ASP	2.2
1	D	180	LEU	2.2
1	D	90	ALA	2.2
1	A	77	SER	2.2
1	C	111	VAL	2.2
1	D	76	ALA	2.1
1	A	165	TYR	2.1
1	C	49	TYR	2.1
1	B	98	GLY	2.1
1	A	175	ILE	2.1
1	D	26	ALA	2.1
1	C	3	GLU	2.1
1	C	43	LYS	2.1
1	A	123	LEU	2.1
1	D	147	TYR	2.1
1	A	201	GLY	2.1
1	A	47	ASP	2.1
1	C	101	ASP	2.1
1	B	177	SER	2.1
1	D	34	PHE	2.1
1	C	183	ALA	2.1
1	D	199	GLN	2.1
1	A	172	SER	2.1
1	B	29	GLU	2.1
1	D	115	GLU	2.1
1	A	27	GLY	2.1
1	B	71	ILE	2.1
1	C	166	TYR	2.1
1	C	122	ALA	2.0
1	D	25	ALA	2.0
1	D	100	ALA	2.0
1	D	122	ALA	2.0
1	D	216	ALA	2.0
1	A	152	LYS	2.0
1	D	193	THR	2.0
1	D	144	GLY	2.0
1	B	167	VAL	2.0
1	D	49	TYR	2.0
1	A	221	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	216	ALA	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](#)

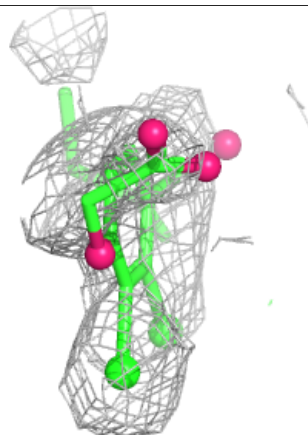
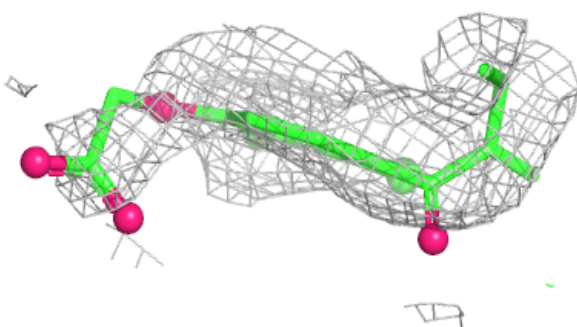
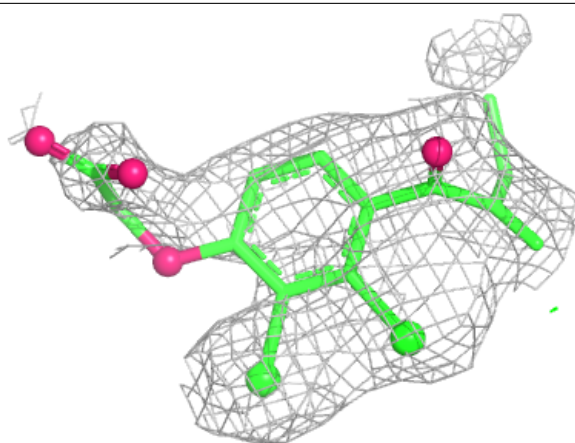
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	EAA	C	223	19/19	0.62	0.17	74,74,74,74	0
2	EAA	A	223	19/19	0.65	0.16	74,74,74,74	0
2	EAA	B	223	19/19	0.70	0.14	74,74,74,74	0
2	EAA	D	223	19/19	0.73	0.15	74,74,74,74	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.

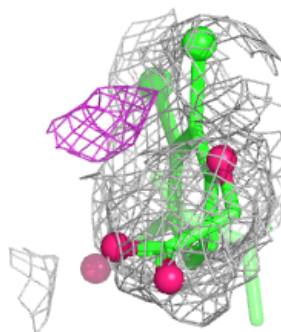
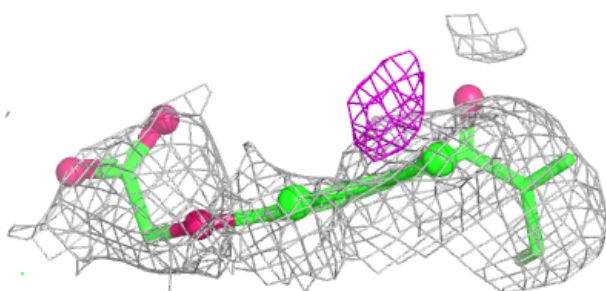
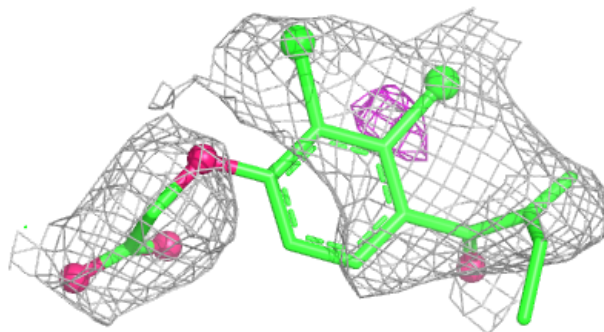
Electron density around EAA C 223:

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

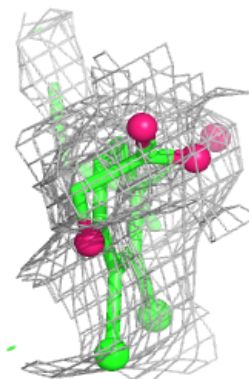
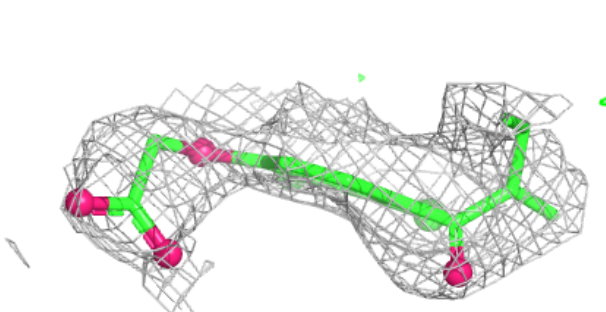
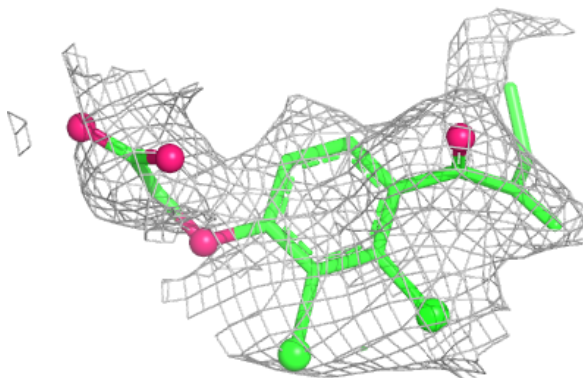


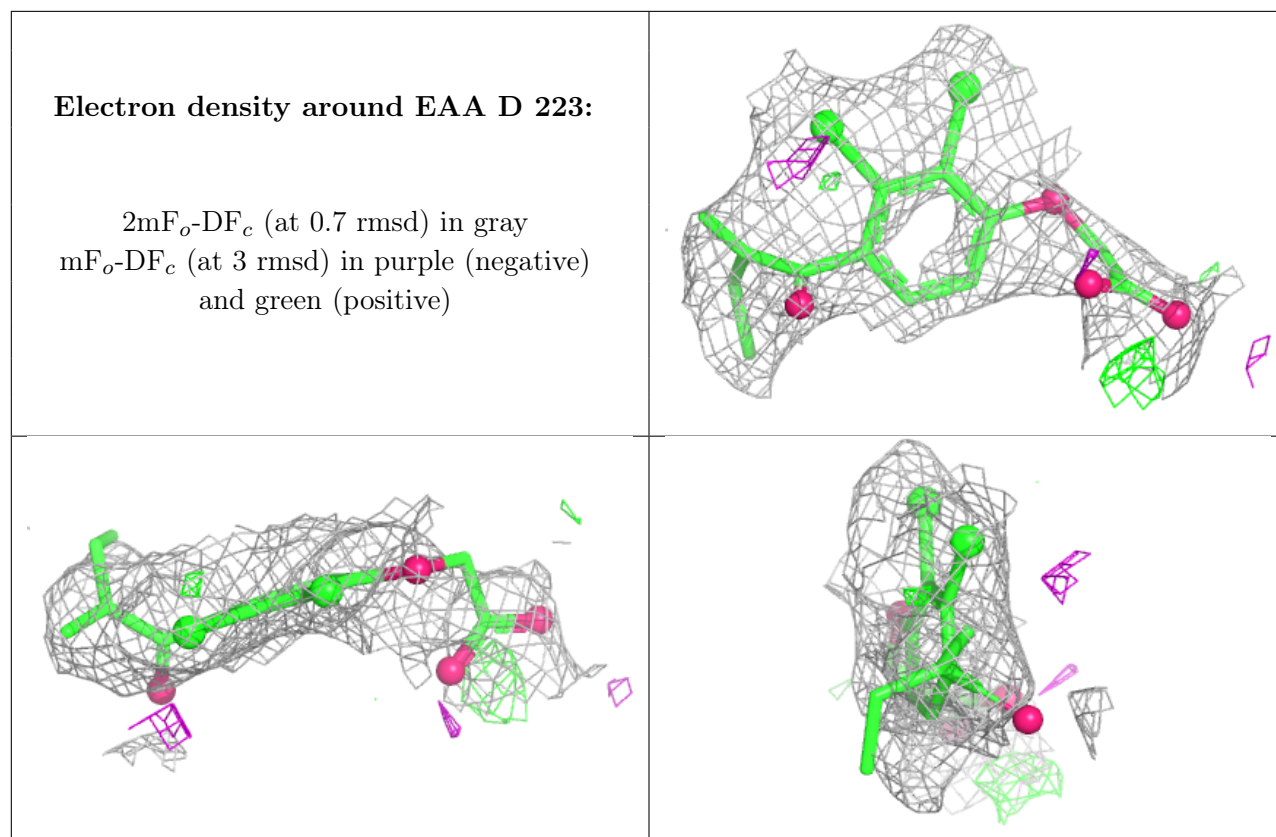
Electron density around EAA A 223:

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

**Electron density around EAA B 223:**

$2mF_o-DF_c$ (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.