



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 10:49 PM UTC

PDB ID : 8OXW / pdb_00008oxw
Title : Transglutaminase 3 in complex with DH patient-derived Fab DH63-B02
Authors : Heggelund, J.E.; Sollid, L.M.
Deposited on : 2023-05-02
Resolution : 1.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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
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

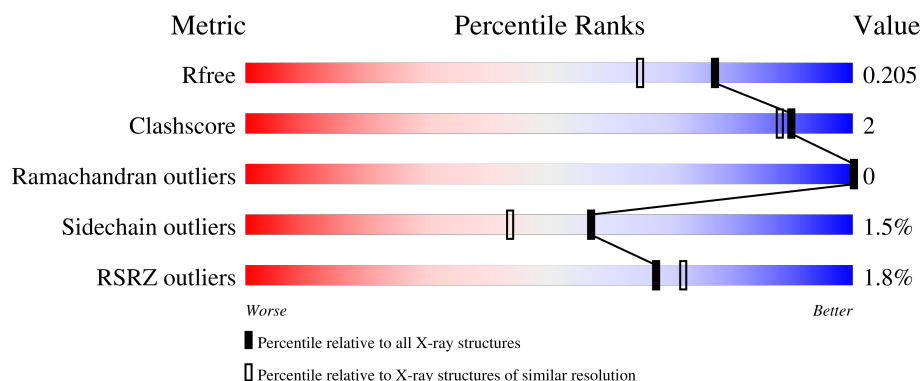
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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))
R_{free}	180053	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	693	2% 92% 6%
2	B	225	2% 93% 6%
3	C	216	% 92% 7%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 18576 atoms, of which 8640 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 Protein-glutamine gamma-glutamyltransferase E 27 kDa non-catalytic chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	679	Total	C	H	N	O	S	139	10	0
			10687	3386	5314	933	1028	26			

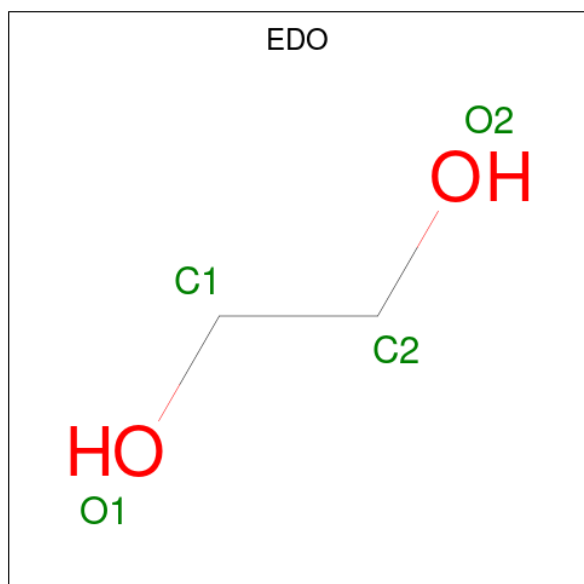
- Molecule 2 is a protein called Antibody fab fragment heavy chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	225	Total	C	H	N	O	S	73	4	0
			3412	1079	1687	295	342	9			

- Molecule 3 is a protein called Antibody fab fragment light chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	216	Total	C	H	N	O	S	74	1	0
			3208	1013	1573	274	342	6			

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C H O 10 2 6 2	1	0
4	A	1	Total C H O 10 2 6 2	1	0
4	A	1	Total C H O 10 2 6 2	1	0
4	A	1	Total C H O 10 2 6 2	1	0
4	A	1	Total C H O 10 2 6 2	1	0
4	A	1	Total C H O 10 2 6 2	1	0
4	A	1	Total C H O 10 2 6 2	1	0
4	A	1	Total C H O 10 2 6 2	1	0
4	A	1	Total C H O 10 2 6 2	1	0
4	A	1	Total C H O 10 2 6 2	1	0
4	C	1	Total C H O 10 2 6 2	1	0

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	3	Total Ca 3 3	0	0

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total Cl 1 1	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	631	Total O 632 632	0	1
7	B	230	Total O 231 231	0	1

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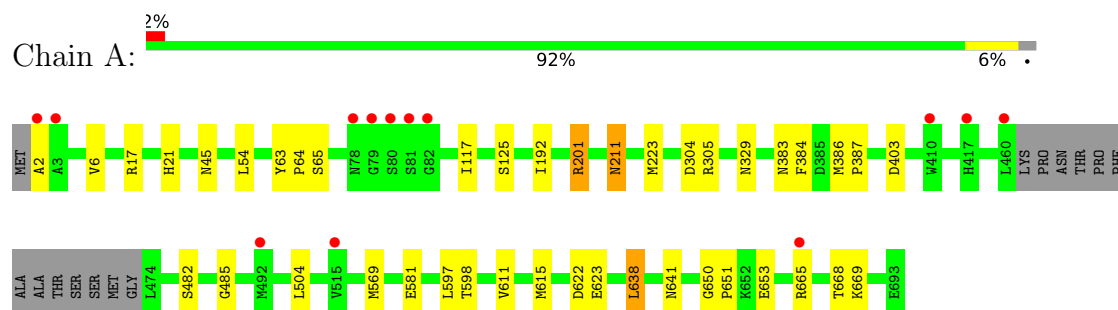
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	290	Total 292	O 292	0	2

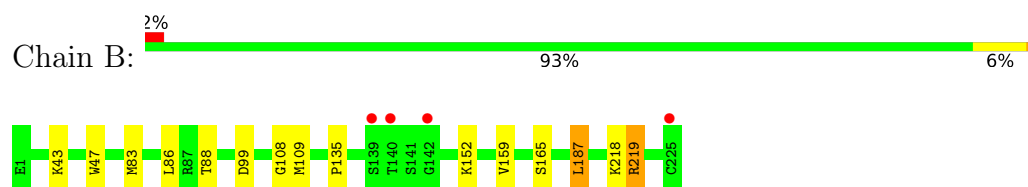
3 Residue-property plots [i](#)

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.

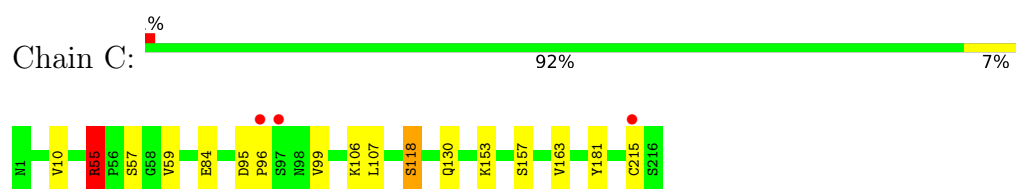
- Molecule 1: Protein-glutamine gamma-glutamyltransferase E 27 kDa non-catalytic chain



- Molecule 2: Antibody fab fragment heavy chain



- Molecule 3: Antibody fab fragment light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.50Å 93.49Å 90.64Å 90.00° 92.63° 90.00°	Depositor
Resolution (Å)	45.31 – 1.70 45.31 – 1.70	Depositor EDS
% Data completeness (in resolution range)	98.4 (45.31-1.70) 97.1 (45.31-1.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.17 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.8.0405	Depositor
R, R_{free}	0.172 , 0.210 0.166 , 0.205	Depositor DCC
R_{free} test set	7482 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	26.2	Xtriage
Anisotropy	0.516	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 43.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.016 for -h,-l,-k 0.000 for -h,l,k 0.025 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	18576	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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: CA, EDO, CL

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	0.76	1/5490 (0.0%)	1.13	6/7443 (0.1%)
2	B	0.75	0/1768	1.12	4/2403 (0.2%)
3	C	0.80	1/1674 (0.1%)	1.19	2/2284 (0.1%)
All	All	0.77	2/8932 (0.0%)	1.14	12/12130 (0.1%)

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	A	0	2
2	B	0	1
3	C	0	1
All	All	0	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	84	GLU	CD-OE1	5.76	1.36	1.25
1	A	21	HIS	ND1-CE1	5.53	1.38	1.32

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	641	ASN	CA-CB-CG	7.02	119.62	112.60
1	A	211	ASN	CA-CB-CG	6.55	119.15	112.60
1	A	403	ASP	CA-CB-CG	6.53	119.13	112.60
3	C	118	SER	CB-CA-C	6.21	119.78	109.53
2	B	152	LYS	N-CA-CB	-5.49	102.06	110.85

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	17	ARG	Sidechain
1	A	201[A]	ARG	Sidechain
2	B	219	ARG	Sidechain
3	C	55	ARG	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	5373	5314	5298	17	0
2	B	1725	1687	1682	7	0
3	C	1635	1573	1568	9	0
4	A	40	60	60	0	0
4	C	4	6	6	0	0
5	A	3	0	0	0	0
6	A	1	0	0	0	0
7	A	632	0	0	1	0
7	B	231	0	0	1	0
7	C	292	0	0	1	0
All	All	9936	8640	8614	32	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 2.

The worst 5 of 32 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:201[A]:ARG:NH1	7:A:801:HOH:O	1.98	0.95
1:A:54:LEU:CD2	1:A:117:ILE:HG12	2.35	0.56
3:C:95:ASP:HB2	3:C:96:PRO:CD	2.36	0.56
1:A:622:ASP:C	1:A:651:PRO:HB3	2.31	0.56
3:C:106[A]:LYS:NZ	7:C:404:HOH:O	2.38	0.55

There are no symmetry-related clashes.

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	685/693 (99%)	662 (97%)	23 (3%)	0	100	100
2	B	227/225 (101%)	221 (97%)	6 (3%)	0	100	100
3	C	215/216 (100%)	206 (96%)	9 (4%)	0	100	100
All	All	1127/1134 (99%)	1089 (97%)	38 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	594/595 (100%)	587 (99%)	7 (1%)	63	51
2	B	195/191 (102%)	191 (98%)	4 (2%)	47	30
3	C	190/189 (100%)	187 (98%)	3 (2%)	55	41
All	All	979/975 (100%)	965 (99%)	14 (1%)	57	46

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	109	MET
2	B	159	VAL
3	C	215	CYS
3	C	57	SER
3	C	118	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	500	ASN
2	B	39	GLN
3	C	39	GLN
2	B	82	GLN
1	A	89	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 4 are monoatomic - leaving 11 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	$\# Z > 2$
4	EDO	A	701	-	3,3,3	0.23	0	2,2,2	0.58	0
4	EDO	A	702	-	3,3,3	0.24	0	2,2,2	0.55	0
4	EDO	A	707	-	3,3,3	0.48	0	2,2,2	0.68	0
4	EDO	A	708	-	3,3,3	0.48	0	2,2,2	0.37	0
4	EDO	A	705	-	3,3,3	0.06	0	2,2,2	0.11	0
4	EDO	A	710	-	3,3,3	0.65	0	2,2,2	0.66	0
4	EDO	C	301	-	3,3,3	0.31	0	2,2,2	0.39	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	A	706	-	3,3,3	0.23	0	2,2,2	0.43	0
4	EDO	A	704	-	3,3,3	0.43	0	2,2,2	0.65	0
4	EDO	A	703	-	3,3,3	0.28	0	2,2,2	0.05	0
4	EDO	A	709	-	3,3,3	0.23	0	2,2,2	0.41	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	EDO	A	701	-	-	1/1/1/1	-
4	EDO	A	702	-	-	0/1/1/1	-
4	EDO	A	707	-	-	1/1/1/1	-
4	EDO	A	708	-	-	0/1/1/1	-
4	EDO	A	705	-	-	1/1/1/1	-
4	EDO	A	710	-	-	1/1/1/1	-
4	EDO	C	301	-	-	1/1/1/1	-
4	EDO	A	706	-	-	0/1/1/1	-
4	EDO	A	704	-	-	1/1/1/1	-
4	EDO	A	703	-	-	0/1/1/1	-
4	EDO	A	709	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	701	EDO	O1-C1-C2-O2
4	A	705	EDO	O1-C1-C2-O2
4	A	704	EDO	O1-C1-C2-O2
4	A	710	EDO	O1-C1-C2-O2
4	A	707	EDO	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	679/693 (97%)	-0.10	13 (1%) 66 71	9, 33, 64, 99	10 (1%)
2	B	225/225 (100%)	-0.10	4 (1%) 67 72	19, 33, 64, 115	4 (1%)
3	C	216/216 (100%)	-0.36	3 (1%) 73 77	13, 29, 49, 154	1 (0%)
All	All	1120/1134 (98%)	-0.15	20 (1%) 67 72	9, 33, 63, 154	15 (1%)

The worst 5 of 20 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	215	CYS	5.6
1	A	460	LEU	4.3
1	A	79	GLY	3.2
2	B	140	THR	3.2
3	C	96	PRO	2.9

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](#)

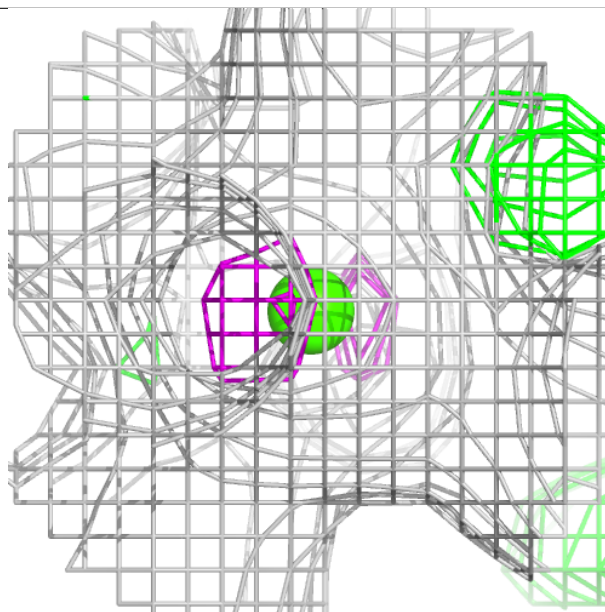
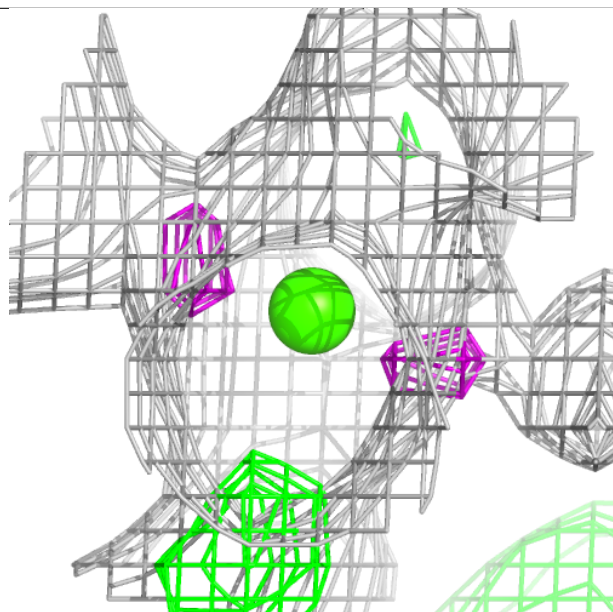
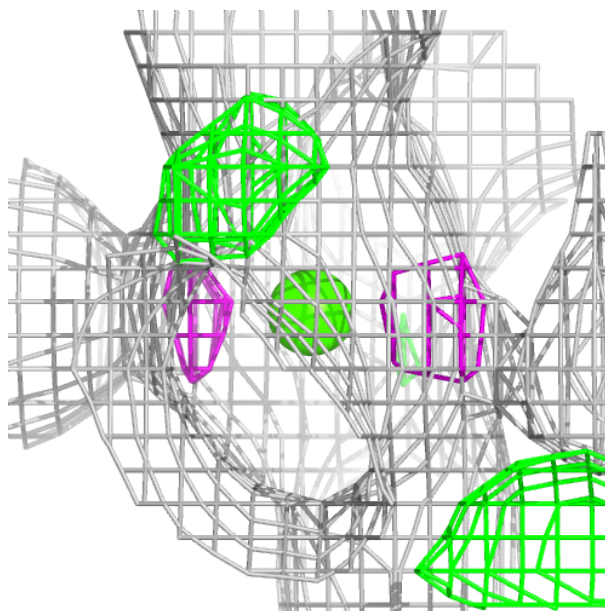
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists 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
4	EDO	A	709	4/4	0.79	0.14	65,73,74,75	1
4	EDO	A	706	4/4	0.88	0.12	57,60,63,65	1
4	EDO	A	703	4/4	0.88	0.12	48,48,59,59	1
4	EDO	A	707	4/4	0.90	0.12	47,56,59,61	1
4	EDO	A	701	4/4	0.90	0.11	43,54,55,55	1
4	EDO	A	708	4/4	0.91	0.11	45,48,55,55	1
4	EDO	A	705	4/4	0.93	0.11	60,63,64,64	1
4	EDO	A	710	4/4	0.93	0.11	44,58,59,65	1
4	EDO	C	301	4/4	0.94	0.10	43,46,55,55	1
4	EDO	A	704	4/4	0.95	0.13	29,44,72,74	1
4	EDO	A	702	4/4	0.97	0.05	31,35,35,36	1
5	CA	A	713	1/1	0.98	0.05	37,37,37,37	0
6	CL	A	714	1/1	0.98	0.09	46,46,46,46	0
5	CA	A	711	1/1	0.99	0.02	22,22,22,22	0
5	CA	A	712	1/1	0.99	0.04	35,35,35,35	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.

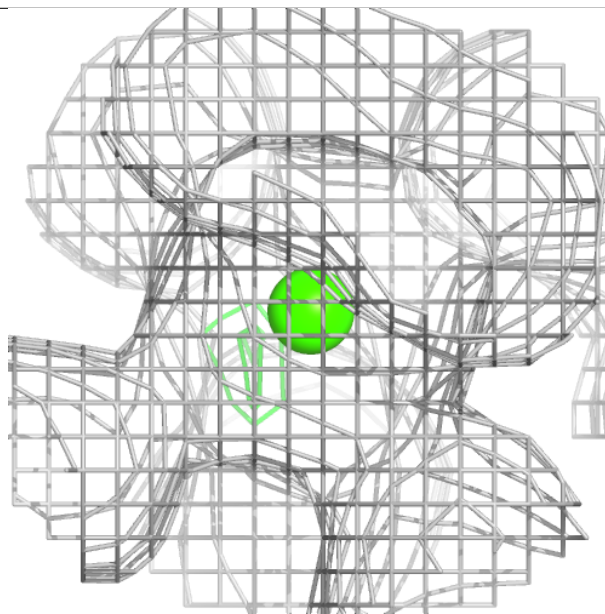
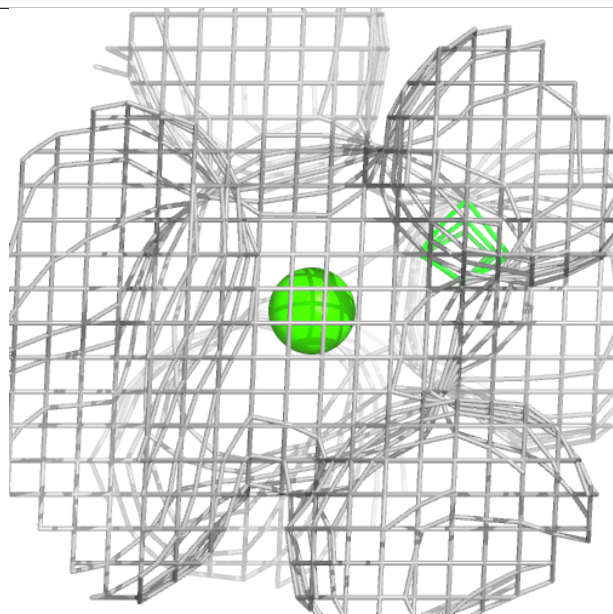
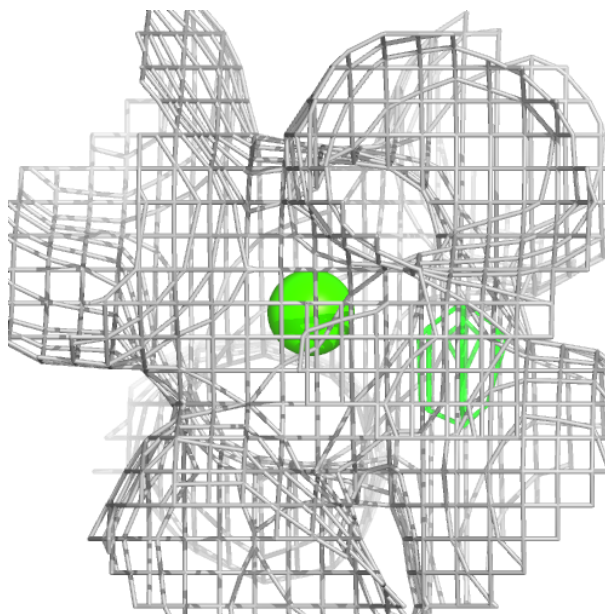
Electron density around CA A 713:

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



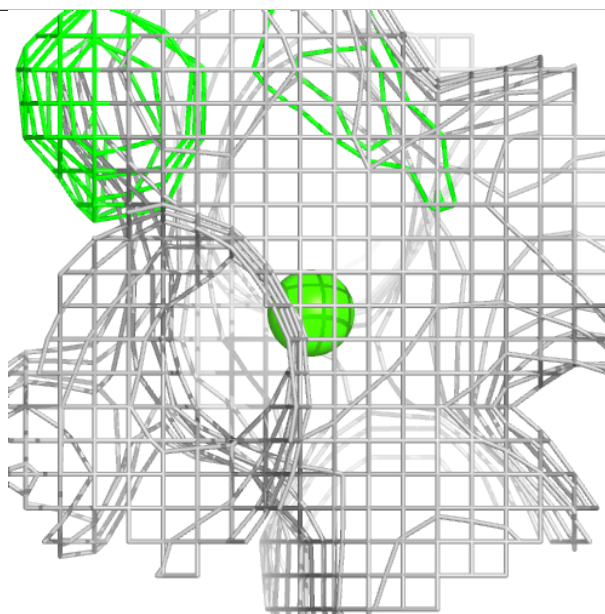
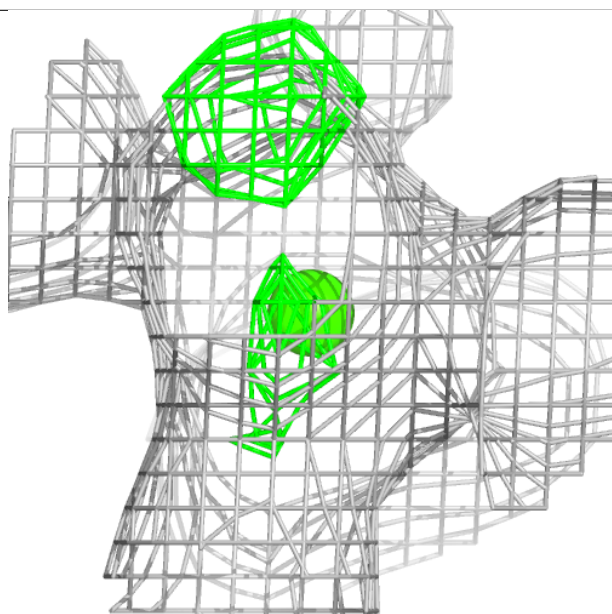
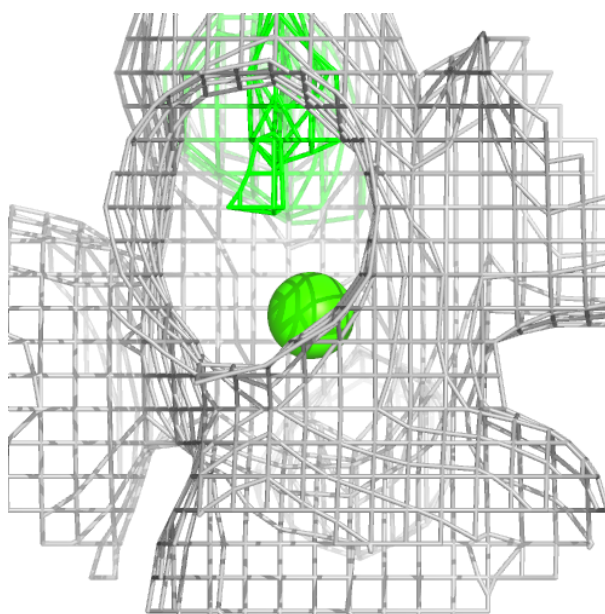
Electron density around CA A 711:

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



Electron density around CA A 712:

$2mF_o-DF_c$ (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.