



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 02:31 PM EDT

PDB ID : 5E2N / pdb_00005e2n
Title : Crystal structure of human carbonic anhydrase isozyme XIII with 3-(cyclooctylamino)-2,5,6-trifluoro-4-[(2-hydroxyethyl)sulfonyl]benzenesulfonamide
Authors : Manakova, E.; Smirnov, A.; Grazulis, S.
Deposited on : 2015-10-01
Resolution : 1.53 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

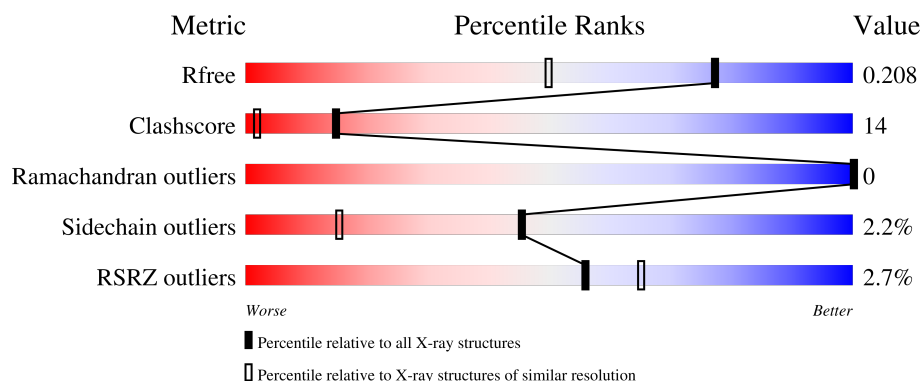
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.53 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	1003 (1.54-1.54)
Clashscore	190562	1025 (1.54-1.54)
Ramachandran outliers	187476	1007 (1.54-1.54)
Sidechain outliers	187428	1007 (1.54-1.54)
RSRZ outliers	180081	1002 (1.54-1.54)

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	263	
1	B	263	

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
3	PEG	A	302	-	X	X	-
5	V14	A	307	-	-	X	-
5	V14	B	306	-	-	X	-
6	CIT	B	302	-	X	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 4795 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 Carbonic anhydrase 13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	10	0
			2119	1351	362	405	1			
1	B	258	Total	C	N	O	S	0	6	0
			2087	1327	364	395	1			

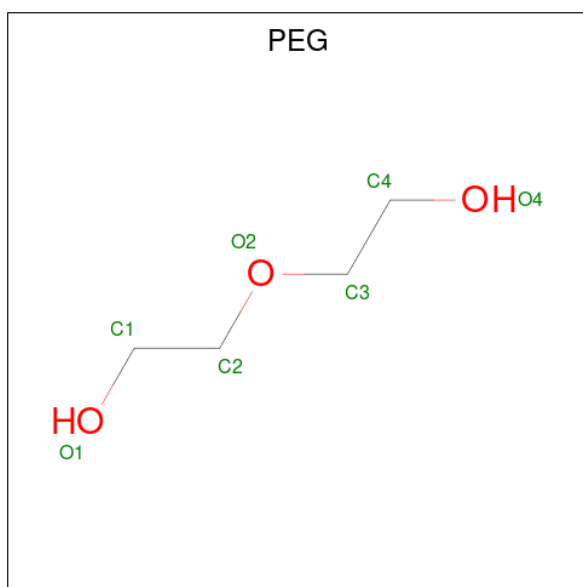
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP Q8N1Q1
B	1	MET	-	initiating methionine	UNP Q8N1Q1

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

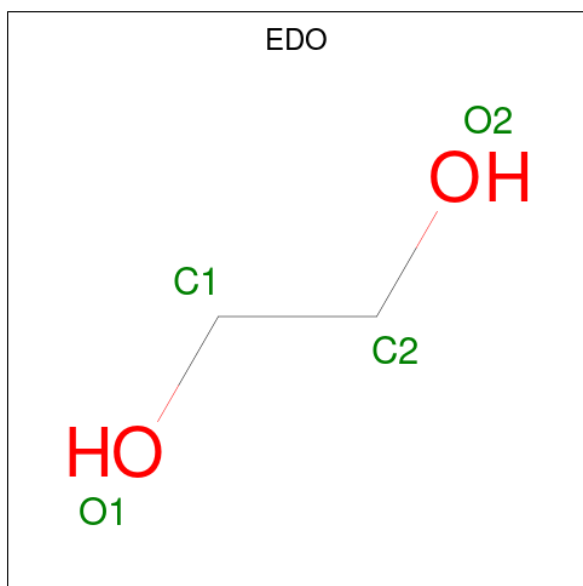
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



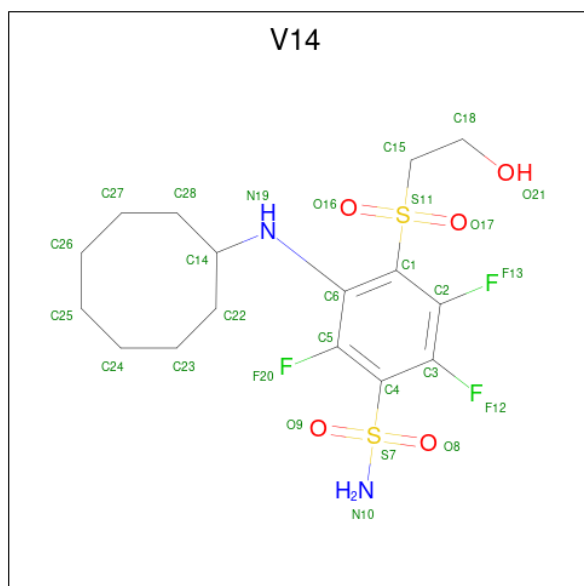
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		

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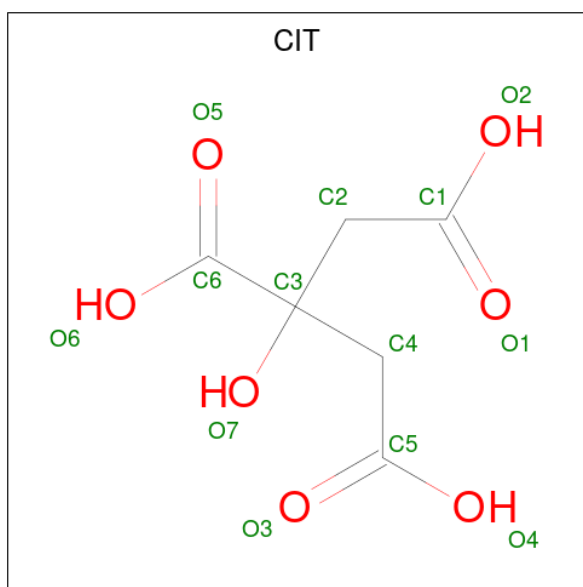
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is 3-(cyclooctylamino)-2,5,6-trifluoro-4-[(2-hydroxyethyl)sulfonyl]benzenesulfonamide (CCD ID: V14) (formula: $C_{16}H_{23}F_3N_2O_5S_2$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	C	F	N	O	S	0	0
			28	16	3	2	5	2		
5	B	1	Total	C	F	N	O	S	0	0
			28	16	3	2	5	2		

- Molecule 6 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			13	6	7		

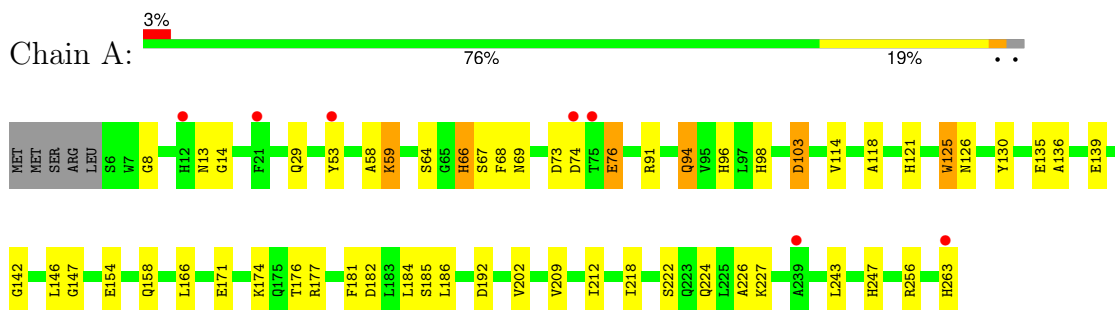
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	224	Total	O	0	0
			224	224		
7	B	259	Total	O	0	0
			259	259		

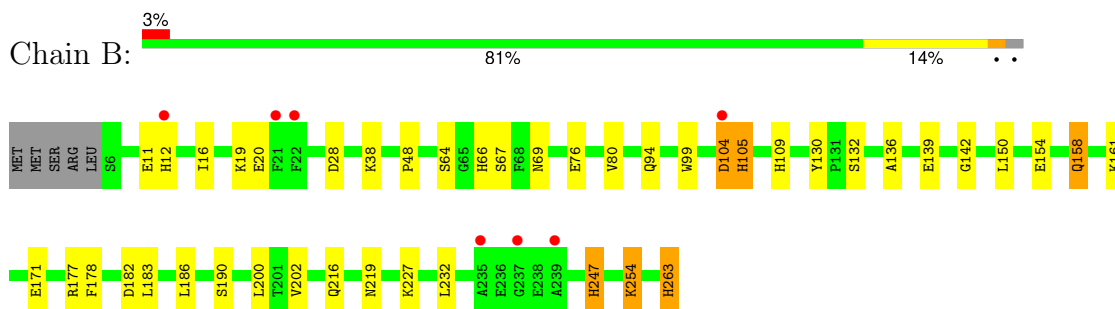
3 Residue-property plots [i](#)

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: Carbonic anhydrase 13



• Molecule 1: Carbonic anhydrase 13



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	56.09Å 57.60Å 159.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	54.18 – 1.53 54.18 – 1.53	Depositor EDS
% Data completeness (in resolution range)	98.7 (54.18-1.53) 98.7 (54.18-1.53)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.76 (at 1.53Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.176 , 0.212 0.173 , 0.208	Depositor DCC
R_{free} test set	7856 reflections (9.94%)	wwPDB-VP
Wilson B-factor (Å ²)	18.2	Xtriage
Anisotropy	0.429	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 39.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.025 for k,h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4795	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CIT, ZN, V14, PEG, EDO

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	1.56	16/2182 (0.7%)	1.42	9/2971 (0.3%)
1	B	1.53	9/2148 (0.4%)	1.32	1/2923 (0.0%)
All	All	1.54	25/4330 (0.6%)	1.37	10/5894 (0.2%)

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	247	HIS	ND1-CE1	10.06	1.42	1.32
1	A	73	ASP	C-O	7.34	1.32	1.24
1	A	96	HIS	CE1-NE2	7.22	1.39	1.32
1	A	243	LEU	CA-C	7.08	1.61	1.53
1	B	109	HIS	ND1-CE1	7.00	1.39	1.32
1	A	13	ASN	N-CA	6.69	1.53	1.46
1	A	98	HIS	ND1-CE1	6.48	1.39	1.32
1	A	73	ASP	CA-C	-6.25	1.44	1.52
1	A	66	HIS	ND1-CE1	6.03	1.38	1.32
1	B	80	VAL	C-O	6.02	1.30	1.23
1	A	126	ASN	C-O	-5.85	1.16	1.24
1	B	254	LYS	CA-C	5.70	1.60	1.53
1	A	118	ALA	N-CA	5.62	1.52	1.46
1	A	247	HIS	ND1-CE1	5.55	1.38	1.32
1	B	183	LEU	N-CA	-5.53	1.39	1.46
1	B	12	HIS	CG-CD2	5.43	1.41	1.35
1	A	243	LEU	CA-CB	5.33	1.59	1.52
1	B	150	LEU	N-CA	5.30	1.52	1.46
1	A	130	TYR	CA-CB	5.29	1.61	1.53
1	B	263	HIS	CG-CD2	5.29	1.41	1.35
1	A	66	HIS	CG-CD2	5.29	1.41	1.35
1	A	76	GLU	N-CA	-5.29	1.39	1.45
1	A	121	HIS	CG-CD2	5.23	1.41	1.35
1	A	125	TRP	CA-C	5.21	1.58	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	105	HIS	ND1-CE1	5.08	1.37	1.32

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	ASP	N-CA-C	6.28	120.04	112.38
1	A	209	VAL	N-CA-C	5.88	116.40	108.17
1	A	73	ASP	N-CA-C	-5.63	101.11	109.56
1	A	68	PHE	N-CA-C	-5.51	101.55	110.32
1	A	114	VAL	N-CA-C	5.47	116.64	108.54
1	A	166	LEU	N-CA-C	5.47	117.24	111.28
1	A	94	GLN	N-CA-C	5.43	117.10	108.79
1	B	177	ARG	CA-CB-CG	-5.36	103.39	114.10
1	A	226	ALA	N-CA-C	5.11	118.29	111.75
1	A	147	GLY	CA-C-O	-5.03	115.71	120.64

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	2119	0	2023	48	0
1	B	2087	0	2011	52	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	7	0	9	4	0
4	A	16	0	24	5	0
4	B	12	0	18	2	0
5	A	28	0	23	24	0
5	B	28	0	23	24	0
6	B	13	0	4	2	0
7	A	224	0	0	6	1
7	B	259	0	0	11	1
All	All	4795	0	4135	117	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 14.

All (117) 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)
5:A:307:V14:H18	5:A:307:V14:N19	1.45	1.17
1:A:69:ASN:HD21	5:A:307:V14:C25	1.66	1.07
5:A:307:V14:H8	5:A:307:V14:C24	1.72	1.02
1:A:69:ASN:HD21	5:A:307:V14:H16	1.26	0.99
1:A:76:GLU:HG3	7:A:405:HOH:O	1.68	0.93
1:A:69:ASN:ND2	5:A:307:V14:H16	1.82	0.93
5:A:307:V14:C24	5:A:307:V14:O17	2.17	0.93
1:B:158[A]:GLN:NE2	7:B:401:HOH:O	2.02	0.91
1:B:202:VAL:HG21	5:B:306:V14:H9	1.55	0.89
5:A:307:V14:H18	5:A:307:V14:O17	1.73	0.88
1:A:69:ASN:HD21	5:A:307:V14:H17	1.43	0.83
1:A:69:ASN:ND2	5:A:307:V14:C25	2.40	0.83
5:A:307:V14:O17	5:A:307:V14:H19	1.82	0.79
1:A:256:ARG:HG3	3:A:302:PEG:H11	1.69	0.75
1:A:69:ASN:CG	5:A:307:V14:H16	2.11	0.75
1:B:104:ASP:CG	7:B:402:HOH:O	2.31	0.74
1:A:176[A]:THR:CG2	7:A:491:HOH:O	2.36	0.72
1:A:202:VAL:HG21	5:A:307:V14:H9	1.70	0.71
5:A:307:V14:N19	5:A:307:V14:C24	2.39	0.71
1:B:202:VAL:HG21	5:B:306:V14:C14	2.20	0.71
1:A:224:GLN:HA	1:A:227:LYS:HE2	1.73	0.70
5:B:306:V14:H20	5:B:306:V14:C18	2.22	0.70
1:A:69:ASN:OD1	5:A:307:V14:H16	1.91	0.69
1:B:66:HIS:CD2	5:B:306:V14:H19	2.26	0.69
1:B:66:HIS:HD2	5:B:306:V14:C24	2.06	0.68
5:A:307:V14:N19	5:A:307:V14:O17	2.24	0.68
1:B:19[B]:LYS:HG2	1:B:20:GLU:N	2.10	0.66
1:B:202:VAL:HG21	5:B:306:V14:H23	1.78	0.66
1:B:64:SER:CB	5:B:306:V14:H17	2.26	0.65
1:A:154:GLU:HB3	7:A:567:HOH:O	1.96	0.65
1:B:19[A]:LYS:HE2	7:B:450:HOH:O	1.97	0.65
5:B:306:V14:O21	5:B:306:V14:H16	1.98	0.64
1:B:16:ILE:O	1:B:19[A]:LYS:HE3	1.99	0.63
5:B:306:V14:H20	5:B:306:V14:H5	1.79	0.63
5:B:306:V14:H9	5:B:306:V14:F20	1.88	0.63
4:A:303:EDO:H21	7:B:440:HOH:O	1.97	0.63
1:B:66:HIS:HD2	5:B:306:V14:H19	1.63	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:ALA:O	1:A:59:LYS:HE3	2.03	0.59
1:A:177[B]:ARG:NH1	7:A:403:HOH:O	2.35	0.59
1:A:182:ASP:HB3	4:A:304:EDO:H11	1.84	0.59
1:B:104:ASP:OD2	7:B:402:HOH:O	2.16	0.58
1:A:146:LEU:HD12	1:A:212:ILE:HB	1.84	0.58
1:A:67:SER:HB3	5:A:307:V14:H13	1.86	0.58
1:B:158[B]:GLN:HG3	1:B:186:LEU:CD1	2.33	0.58
6:B:302:CIT:O3	6:B:302:CIT:C6	2.50	0.57
5:B:306:V14:C18	5:B:306:V14:H8	2.19	0.55
1:B:158[B]:GLN:HG2	7:B:555:HOH:O	2.07	0.55
1:B:190:SER:H	1:B:216:GLN:NE2	2.05	0.54
1:B:202:VAL:HG21	5:B:306:V14:C22	2.36	0.54
1:B:99:TRP:O	1:B:247:HIS:HD2	1.90	0.54
1:A:29:GLN:OE1	3:A:302:PEG:H22	2.07	0.54
1:B:254:LYS:HE2	7:B:444:HOH:O	2.07	0.53
1:B:19[B]:LYS:O	1:B:20:GLU:C	2.49	0.53
1:A:29:GLN:HG2	3:A:302:PEG:H12	1.90	0.53
1:A:69:ASN:ND2	5:A:307:V14:H17	2.15	0.53
4:A:303:EDO:H11	1:B:48:PRO:HD3	1.92	0.52
1:B:158[B]:GLN:HG3	1:B:186:LEU:HD13	1.92	0.51
1:A:135[B]:GLU:OE2	1:A:139[B]:GLU:OE2	2.27	0.51
1:B:132:SER:HA	4:B:304:EDO:H12	1.92	0.51
1:A:74[B]:ASP:CG	1:A:125:TRP:HE1	2.18	0.51
1:B:130:TYR:CE1	1:B:139:GLU:HG3	2.46	0.51
1:A:64:SER:CB	5:A:307:V14:H17	2.41	0.50
1:B:104:ASP:OD1	7:B:402:HOH:O	2.20	0.50
1:A:64:SER:OG	5:A:307:V14:H15	2.11	0.50
5:A:307:V14:H18	5:A:307:V14:H8	0.76	0.50
1:B:69:ASN:HD21	5:B:306:V14:H17	1.75	0.49
1:B:202:VAL:CG2	5:B:306:V14:H9	2.37	0.49
1:B:66:HIS:CD2	5:B:306:V14:C24	2.88	0.49
1:B:136:ALA:O	1:B:142:GLY:HA3	2.12	0.49
1:B:28:ASP:HA	1:B:254:LYS:HD2	1.95	0.48
1:B:76:GLU:HB2	7:B:478:HOH:O	2.13	0.48
1:B:200:LEU:HD22	5:B:306:V14:C3	2.44	0.48
4:B:304:EDO:H11	7:B:575:HOH:O	2.13	0.48
1:A:74[A]:ASP:C	1:A:91:ARG:HD3	2.39	0.48
1:A:202:VAL:HG21	5:A:307:V14:C14	2.42	0.48
1:A:136:ALA:O	1:A:142:GLY:HA3	2.15	0.47
1:B:105:HIS:HD2	6:B:302:CIT:O5	1.98	0.47
1:B:158[B]:GLN:HG3	1:B:186:LEU:HD12	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158[B]:GLN:HG2	1:B:158[B]:GLN:H	1.54	0.47
1:B:64:SER:HB3	5:B:306:V14:H17	1.97	0.46
1:A:74[A]:ASP:O	1:A:91:ARG:HD2	2.15	0.46
1:B:219[B]:ASN:ND2	1:B:219[B]:ASN:H	2.13	0.46
1:A:64:SER:OG	5:A:307:V14:C26	2.64	0.46
1:B:158[A]:GLN:NE2	1:B:182:ASP:H	2.14	0.45
1:A:74[A]:ASP:O	1:A:91:ARG:CD	2.64	0.45
1:A:224:GLN:HA	1:A:227:LYS:CE	2.45	0.45
1:B:69:ASN:HD21	5:B:306:V14:C25	2.30	0.45
1:B:161:LYS:HG2	1:B:178:PHE:CE1	2.52	0.45
1:B:16:ILE:O	1:B:19[A]:LYS:CE	2.64	0.45
1:B:66:HIS:HD2	5:B:306:V14:H18	1.78	0.45
1:B:227:LYS:HE3	1:B:227:LYS:HB3	1.77	0.45
1:B:67:SER:HB3	5:B:306:V14:H13	1.99	0.44
1:A:135[B]:GLU:CD	1:A:135[B]:GLU:C	2.86	0.44
1:A:53[B]:TYR:HD2	7:A:434:HOH:O	2.00	0.43
1:A:135[B]:GLU:CD	1:A:135[B]:GLU:O	2.61	0.43
1:B:158[B]:GLN:CG	7:B:555:HOH:O	2.66	0.43
1:B:66:HIS:HB2	5:B:306:V14:H19	1.98	0.43
1:A:192:ASP:OD1	4:A:303:EDO:H22	2.19	0.43
5:B:306:V14:C14	5:B:306:V14:F20	2.57	0.43
1:A:186:LEU:O	1:A:218:ILE:HD13	2.19	0.43
1:B:171:GLU:HG3	1:B:232:LEU:HD12	2.01	0.43
1:A:176[A]:THR:HG22	7:A:491:HOH:O	2.14	0.42
1:A:186:LEU:O	1:A:218:ILE:HG21	2.20	0.42
1:A:103:ASP:OD2	1:A:222[B]:SER:OG	2.36	0.42
1:B:69:ASN:OD1	5:B:306:V14:H14	2.19	0.42
1:A:8:GLY:O	1:A:14:GLY:HA2	2.20	0.42
1:B:38:LYS:HB3	1:B:38:LYS:HE2	1.95	0.42
1:A:171:GLU:OE2	1:A:174:LYS:NZ	2.51	0.42
1:A:256:ARG:HG3	3:A:302:PEG:C1	2.43	0.42
1:A:186:LEU:HD23	1:A:186:LEU:HA	1.96	0.42
1:A:67:SER:CB	5:A:307:V14:H13	2.50	0.41
1:A:66:HIS:HD2	5:A:307:V14:H21	1.85	0.41
1:A:181:PHE:CD1	4:A:304:EDO:H22	2.56	0.41
1:A:158:GLN:NE2	1:A:185[B]:SER:OG	2.54	0.40
1:B:190:SER:H	1:B:216:GLN:HE21	1.68	0.40
1:B:219[B]:ASN:H	1:B:219[B]:ASN:HD22	1.70	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)
7:A:410:HOH:O	7:B:457:HOH:O[1_545]	2.13	0.07

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	266/263 (101%)	257 (97%)	9 (3%)	0	100	100
1	B	262/263 (100%)	251 (96%)	11 (4%)	0	100	100
All	All	528/526 (100%)	508 (96%)	20 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	232/232 (100%)	228 (98%)	4 (2%)	53	23
1	B	230/232 (99%)	223 (97%)	7 (3%)	36	8
All	All	462/464 (100%)	451 (98%)	11 (2%)	45	13

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	LYS
1	A	94	GLN
1	A	184	LEU

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Mol	Chain	Res	Type
1	A	263	HIS
1	B	11	GLU
1	B	94	GLN
1	B	104	ASP
1	B	154	GLU
1	B	158[A]	GLN
1	B	158[B]	GLN
1	B	263	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	HIS
1	A	69	ASN
1	A	151	GLN
1	A	158	GLN
1	A	251	GLN
1	B	66	HIS
1	B	105	HIS
1	B	175	GLN
1	B	216	GLN
1	B	247	HIS

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 13 ligands modelled in this entry, 2 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	306	-	3,3,3	0.80	0	2,2,2	1.37	0
5	V14	A	307	-	29,29,29	3.29	8 (27%)	37,43,43	4.65	20 (54%)
4	EDO	B	303	-	3,3,3	0.51	0	2,2,2	0.15	0
4	EDO	B	305	-	3,3,3	0.58	0	2,2,2	0.82	0
4	EDO	B	304	-	3,3,3	0.50	0	2,2,2	0.65	0
4	EDO	A	304	-	3,3,3	0.61	0	2,2,2	0.65	0
4	EDO	A	303	-	3,3,3	0.53	0	2,2,2	1.48	0
4	EDO	A	305	-	3,3,3	0.67	0	2,2,2	1.13	0
6	CIT	B	302	-	12,12,12	2.77	8 (66%)	17,17,17	3.77	14 (82%)
3	PEG	A	302	-	6,6,6	0.99	0	5,5,5	4.35	5 (100%)
5	V14	B	306	2	29,29,29	3.48	8 (27%)	37,43,43	4.70	17 (45%)

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	306	-	-	1/1/1/1	-
5	V14	A	307	-	-	8/20/30/30	0/2/2/2
4	EDO	B	303	-	-	0/1/1/1	-
4	EDO	B	305	-	-	1/1/1/1	-
4	EDO	B	304	-	-	1/1/1/1	-
4	EDO	A	304	-	-	1/1/1/1	-
4	EDO	A	303	-	-	0/1/1/1	-
4	EDO	A	305	-	-	1/1/1/1	-
6	CIT	B	302	-	-	7/16/16/16	-
3	PEG	A	302	-	-	3/4/4/4	-
5	V14	B	306	2	-	14/20/30/30	0/2/2/2

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	306	V14	S7-N10	15.56	1.90	1.60
5	A	307	V14	S7-N10	10.82	1.81	1.60
5	A	307	V14	O17-S11	7.78	1.54	1.44
5	A	307	V14	O16-S11	6.38	1.52	1.44
5	B	306	V14	O16-S11	6.17	1.52	1.44
5	A	307	V14	C22-C14	5.28	1.60	1.52
5	A	307	V14	C4-S7	-5.08	1.70	1.78
6	B	302	CIT	O4-C5	-4.96	1.14	1.30
5	B	306	V14	O17-S11	4.26	1.49	1.44
6	B	302	CIT	C3-C6	-4.10	1.49	1.53
6	B	302	CIT	O7-C3	-3.73	1.36	1.43
5	A	307	V14	O9-S7	-3.38	1.37	1.43
5	B	306	V14	C4-S7	-2.99	1.74	1.78
5	B	306	V14	C22-C14	2.91	1.56	1.52
6	B	302	CIT	C2-C3	-2.61	1.50	1.54
6	B	302	CIT	O6-C6	-2.61	1.21	1.30
5	A	307	V14	F13-C2	2.60	1.38	1.34
5	B	306	V14	C14-N19	2.54	1.51	1.46
5	B	306	V14	C3-C4	-2.52	1.35	1.39
6	B	302	CIT	C2-C1	2.42	1.58	1.50
5	A	307	V14	O8-S7	-2.36	1.39	1.43
6	B	302	CIT	O1-C1	-2.28	1.14	1.22
6	B	302	CIT	O5-C6	-2.25	1.15	1.22
5	B	306	V14	C28-C14	2.23	1.55	1.52

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	306	V14	O8-S7-O9	-17.46	91.94	118.80
5	A	307	V14	C3-C4-C5	14.50	126.59	116.65
5	A	307	V14	O8-S7-O9	-13.94	97.36	118.80
5	A	307	V14	O17-S11-O16	-10.72	106.96	118.45
5	B	306	V14	O17-S11-O16	-9.43	108.35	118.45
5	B	306	V14	O9-S7-N10	-9.36	93.86	107.35
5	B	306	V14	O8-S7-C4	8.02	119.06	107.28
5	B	306	V14	O8-S7-N10	7.24	117.79	107.35
5	B	306	V14	C3-C4-C5	7.03	121.47	116.65
6	B	302	CIT	O7-C3-C6	-6.55	99.67	108.96
5	A	307	V14	O8-S7-N10	6.07	116.10	107.35
6	B	302	CIT	C2-C3-C6	5.98	123.26	110.03
5	B	306	V14	C15-S11-C1	5.96	132.88	108.15
5	A	307	V14	O8-S7-C4	5.96	116.03	107.28
6	B	302	CIT	O6-C6-C3	5.94	124.53	113.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	307	V14	O17-S11-C1	-5.67	92.23	106.59
5	B	306	V14	C4-S7-N10	5.41	118.00	108.26
3	A	302	PEG	C3-O2-C2	5.30	136.47	113.26
3	A	302	PEG	O2-C3-C4	5.01	132.19	110.11
6	B	302	CIT	O7-C3-C4	-4.99	98.00	109.38
6	B	302	CIT	O2-C1-C2	4.61	128.95	114.35
5	A	307	V14	C4-S7-N10	4.59	116.52	108.26
3	A	302	PEG	O4-C4-C3	4.33	137.30	111.82
5	A	307	V14	C15-S11-C1	4.23	125.68	108.15
6	B	302	CIT	O7-C3-C2	-4.07	100.09	109.38
5	A	307	V14	O17-S11-C15	3.94	114.63	108.18
5	B	306	V14	O16-S11-C1	-3.83	96.88	106.59
5	B	306	V14	C2-C3-C4	-3.78	117.21	121.78
5	A	307	V14	C6-C5-C4	-3.65	115.56	121.81
5	A	307	V14	O9-S7-C4	-3.53	102.09	107.28
3	A	302	PEG	O1-C1-C2	3.50	132.42	111.82
5	B	306	V14	F12-C3-C2	3.46	126.13	119.27
5	A	307	V14	C2-C3-C4	-3.42	117.64	121.78
6	B	302	CIT	O1-C1-C2	-3.32	113.56	122.95
6	B	302	CIT	C3-C2-C1	-3.29	104.94	113.92
3	A	302	PEG	O2-C2-C1	3.25	124.43	110.11
6	B	302	CIT	C4-C3-C2	3.24	117.62	109.31
5	B	306	V14	O21-C18-C15	-3.19	102.16	109.82
6	B	302	CIT	O4-C5-O3	-3.08	115.40	123.33
5	B	306	V14	F13-C2-C3	-3.05	113.21	119.27
5	A	307	V14	C3-C2-C1	-3.03	118.11	121.78
5	A	307	V14	F12-C3-C2	2.95	125.13	119.27
5	B	306	V14	C2-C1-S11	-2.86	113.86	120.87
5	A	307	V14	C24-C23-C22	-2.81	105.45	117.37
5	A	307	V14	O16-S11-C15	2.79	112.75	108.18
5	A	307	V14	C6-C1-C2	2.73	124.79	119.41
6	B	302	CIT	O2-C1-O1	-2.41	117.14	123.33
5	B	306	V14	F12-C3-C4	-2.28	116.65	120.76
5	B	306	V14	C3-C4-S7	-2.18	116.05	121.34
5	A	307	V14	C2-C1-S11	-2.18	115.53	120.87
5	A	307	V14	C3-C4-S7	-2.16	116.10	121.34
6	B	302	CIT	O5-C6-C3	-2.14	117.94	122.09
6	B	302	CIT	C3-C4-C5	-2.13	108.08	113.92
5	A	307	V14	C23-C24-C25	-2.06	106.96	117.26
5	B	306	V14	F20-C5-C4	-2.02	117.12	120.76
6	B	302	CIT	O4-C5-C4	2.02	120.74	114.35

There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	307	V14	C5-C4-S7-N10
5	A	307	V14	C6-C1-S11-O16
5	A	307	V14	C18-C15-S11-C1
5	A	307	V14	C18-C15-S11-O16
5	A	307	V14	C18-C15-S11-O17
5	A	307	V14	C22-C14-N19-C6
5	B	306	V14	C5-C4-S7-O9
5	B	306	V14	C5-C4-S7-N10
5	B	306	V14	C3-C4-S7-O9
5	B	306	V14	C2-C1-S11-C15
5	B	306	V14	C6-C1-S11-O16
5	B	306	V14	C6-C1-S11-O17
5	B	306	V14	C6-C1-S11-C15
5	B	306	V14	C18-C15-S11-C1
5	B	306	V14	C18-C15-S11-O16
5	B	306	V14	C18-C15-S11-O17
5	B	306	V14	C22-C14-N19-C6
6	B	302	CIT	O7-C3-C4-C5
6	B	302	CIT	C2-C3-C6-O5
6	B	302	CIT	C2-C3-C6-O6
6	B	302	CIT	O7-C3-C6-O5
6	B	302	CIT	O7-C3-C6-O6
6	B	302	CIT	C2-C3-C4-C5
3	A	302	PEG	O2-C3-C4-O4
6	B	302	CIT	C6-C3-C4-C5
4	A	304	EDO	O1-C1-C2-O2
4	B	304	EDO	O1-C1-C2-O2
4	A	306	EDO	O1-C1-C2-O2
3	A	302	PEG	C1-C2-O2-C3
3	A	302	PEG	C4-C3-O2-C2
5	A	307	V14	C28-C14-N19-C6
5	B	306	V14	C28-C14-N19-C6
5	A	307	V14	C2-C1-S11-O16
5	B	306	V14	C3-C4-S7-O8
4	A	305	EDO	O1-C1-C2-O2
4	B	305	EDO	O1-C1-C2-O2
5	B	306	V14	C5-C6-N19-C14

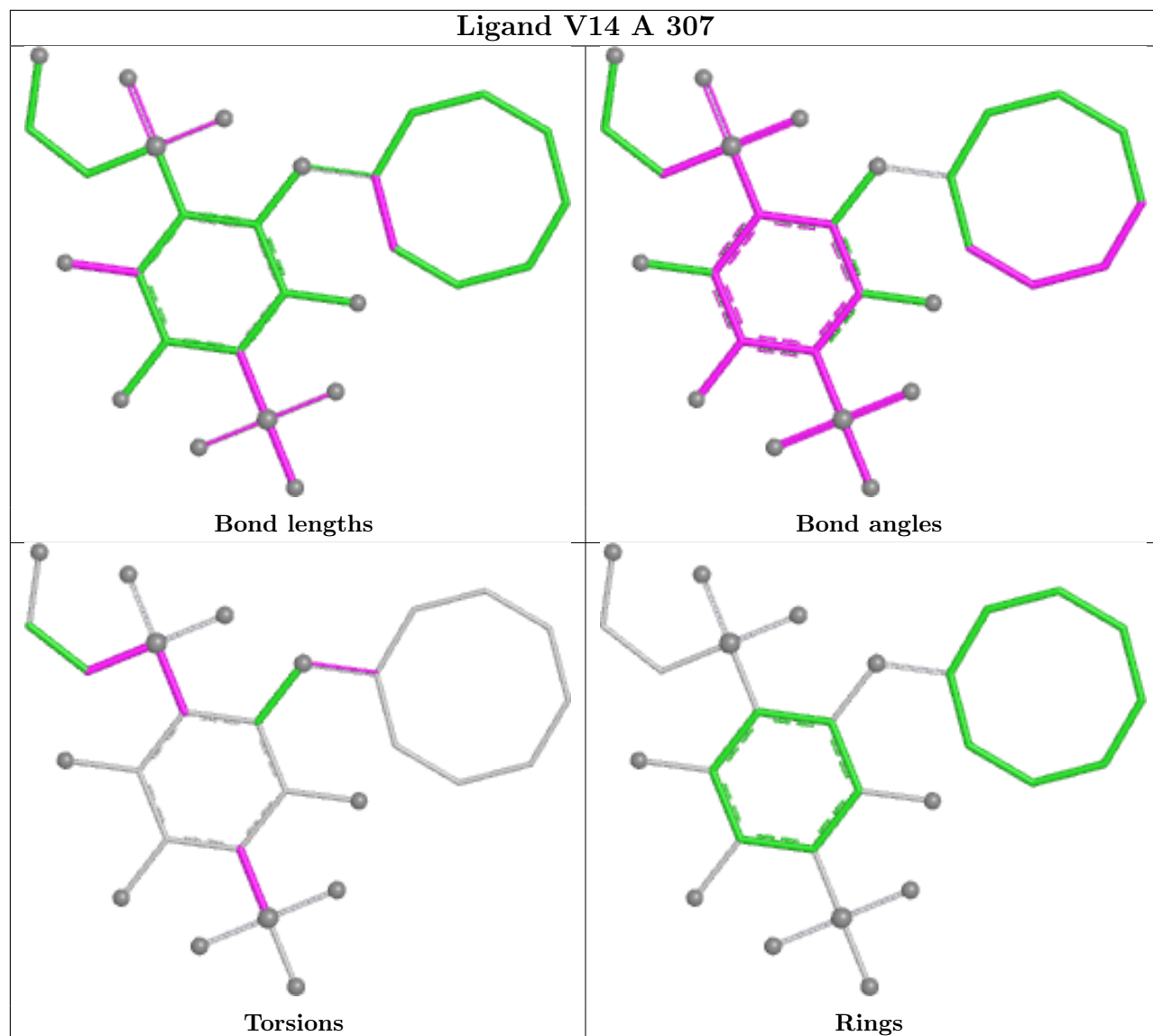
There are no ring outliers.

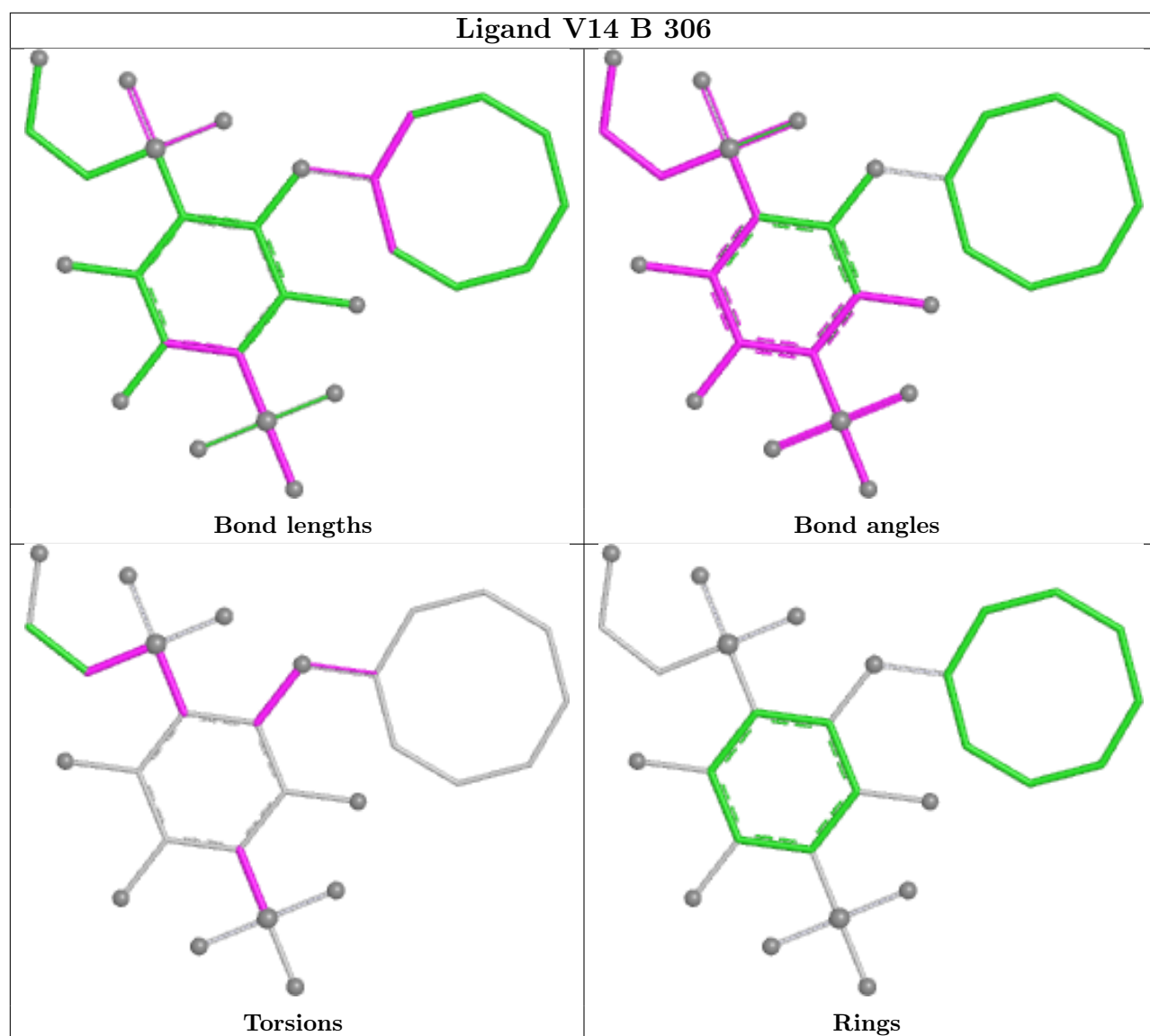
7 monomers are involved in 61 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	307	V14	24	0
4	B	304	EDO	2	0
4	A	304	EDO	2	0
4	A	303	EDO	3	0
6	B	302	CIT	2	0
3	A	302	PEG	4	0
5	B	306	V14	24	0

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

Ligand V14 A 307





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	258/263 (98%)	0.10	7 (2%) 56 65	7, 18, 34, 55	10 (3%)
1	B	258/263 (98%)	-0.01	7 (2%) 56 65	8, 18, 35, 55	6 (2%)
All	All	516/526 (98%)	0.04	14 (2%) 56 65	7, 18, 35, 55	16 (3%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	21	PHE	6.5
1	B	239	ALA	4.1
1	A	12	HIS	3.6
1	A	21	PHE	3.3
1	A	263	HIS	3.2
1	B	237	GLY	3.1
1	A	75	THR	3.0
1	B	104	ASP	2.9
1	A	74[A]	ASP	2.8
1	A	53[A]	TYR	2.6
1	B	235	ALA	2.4
1	B	22	PHE	2.4
1	B	12	HIS	2.2
1	A	239	ALA	2.1

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 ⓘ

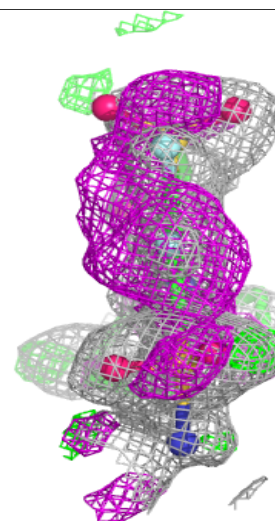
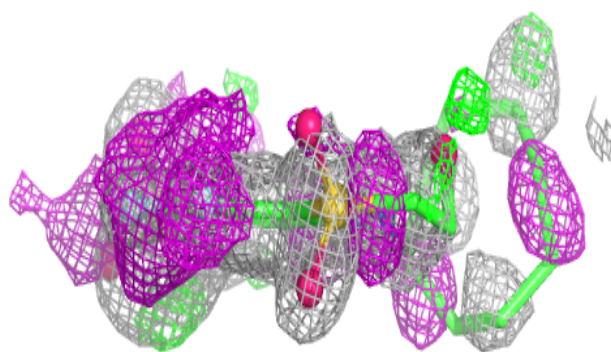
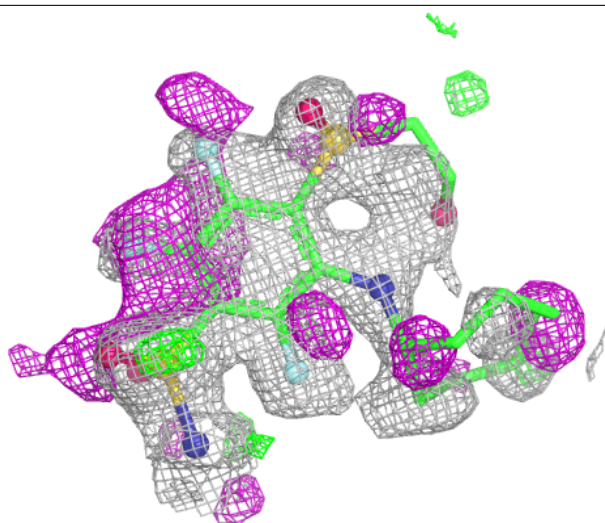
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	EDO	B	305	4/4	0.73	0.15	40,43,43,50	0
6	CIT	B	302	13/13	0.81	0.14	26,48,70,77	0
4	EDO	B	303	4/4	0.82	0.13	31,32,34,38	0
5	V14	B	306	28/28	0.86	0.17	24,39,67,72	0
4	EDO	A	306	4/4	0.86	0.12	26,34,35,35	0
4	EDO	B	304	4/4	0.88	0.12	39,42,44,50	0
4	EDO	A	304	4/4	0.89	0.12	30,34,37,40	0
5	V14	A	307	28/28	0.89	0.15	19,31,57,63	0
4	EDO	A	303	4/4	0.91	0.09	26,29,31,35	0
4	EDO	A	305	4/4	0.93	0.09	26,31,31,32	0
3	PEG	A	302	7/7	0.94	0.08	23,26,29,30	0
2	ZN	B	301	1/1	0.99	0.02	13,13,13,13	0
2	ZN	A	301	1/1	1.00	0.01	11,11,11,11	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.

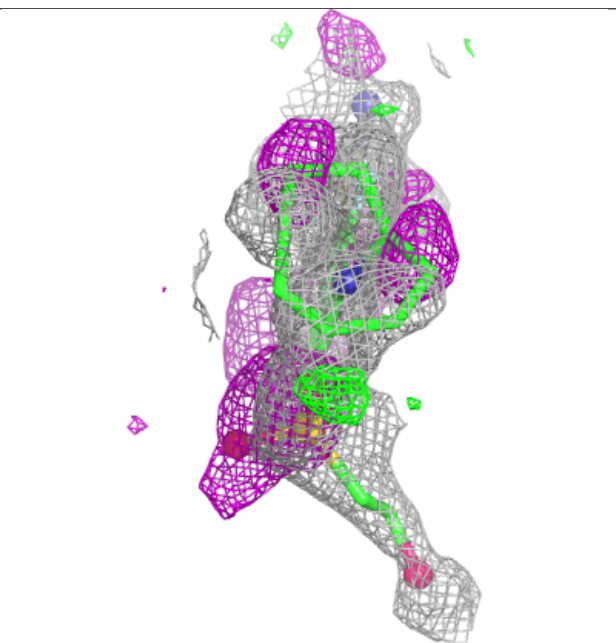
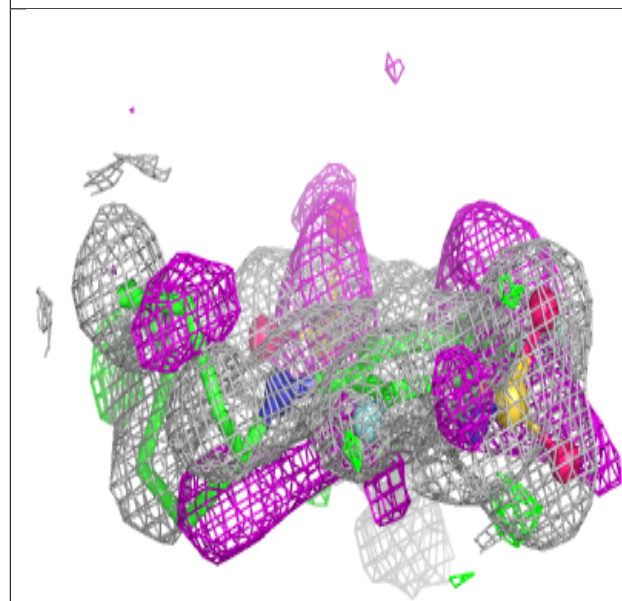
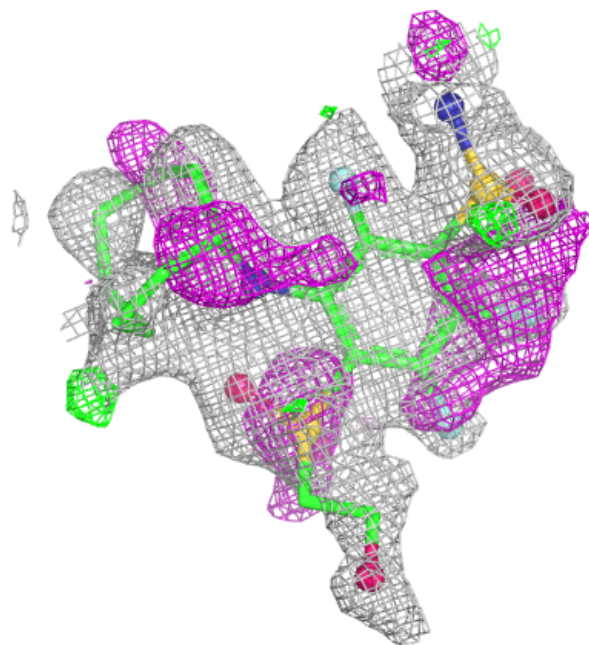
Electron density around V14 B 306:

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



Electron density around V14 A 307:

$2mF_o-DF_c$ (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.