



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:09 PM UTC

PDB ID : 2XO8 / pdb_00002xo8
Title : Crystal Structure of Myosin-2 in Complex with Tribromodichloropseudilin
Authors : Preller, M.; Chinthalapudi, K.; Martin, R.; Knoelker, H.J.; Manstein, D.J.
Deposited on : 2010-08-10
Resolution : 2.40 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

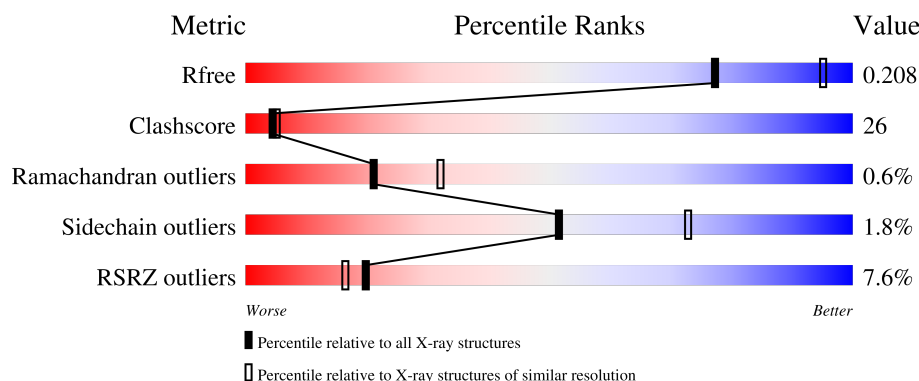
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	776	8% 65% 33% .

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	H70	A	1778	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6720 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.

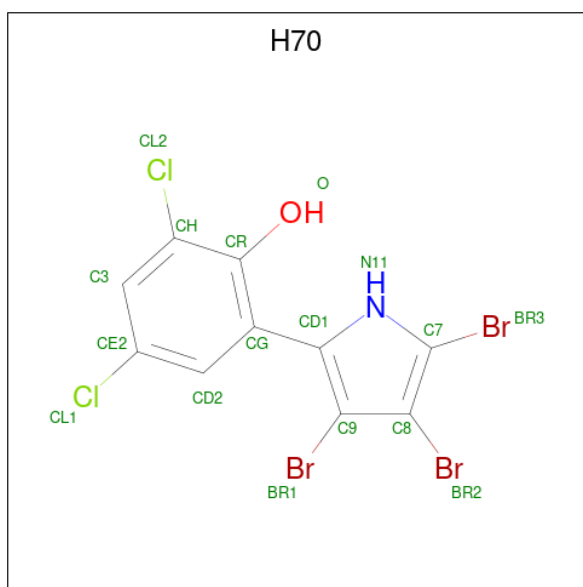
- Molecule 1 is a protein called MYOSIN-2 HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	776	6236	3961	1071	1187	17	36	0	0

There are 19 discrepancies between the modelled and reference sequences:

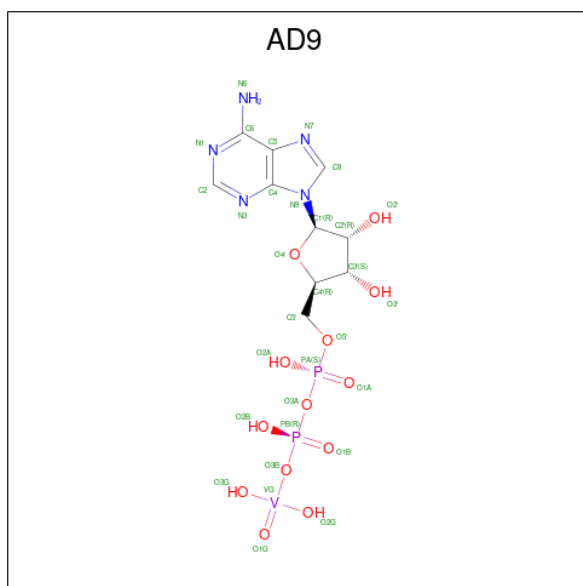
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	ASP	-	expression tag	UNP P08799
A	762	LEU	-	expression tag	UNP P08799
A	763	GLU	-	expression tag	UNP P08799
A	764	SER	-	expression tag	UNP P08799
A	765	ASN	-	expression tag	UNP P08799
A	766	GLU	-	expression tag	UNP P08799
A	767	PRO	-	expression tag	UNP P08799
A	768	PRO	-	expression tag	UNP P08799
A	769	MET	-	expression tag	UNP P08799
A	770	ASP	-	expression tag	UNP P08799
A	771	PHE	-	expression tag	UNP P08799
A	772	ASP	-	expression tag	UNP P08799
A	773	ASP	-	expression tag	UNP P08799
A	774	ASP	-	expression tag	UNP P08799
A	775	ILE	-	expression tag	UNP P08799
A	776	PRO	-	expression tag	UNP P08799
A	777	PHE	-	expression tag	UNP P08799
A	131	ILE	ARG	conflict	UNP P08799
A	435	ILE	LYS	conflict	UNP P08799

- Molecule 2 is 2,4-DICHLORO-6-(3,4,5-TRIBROMO-1H-PYRROL-2-YL)PHENOL (CCD ID: H70) (formula: C₁₀H₄Br₃Cl₂NO).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	Br	C	Cl	N	O		
2	A	1	17	3	10	2	1	1	0	0

- Molecule 3 is ADP METAVANADATE (CCD ID: AD9) (formula: $C_{10}H_{16}N_5O_{13}P_2V$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	N	O	P	V		
3	A	1	31	10	5	13	2	1	0	0

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Mg 1	0	0

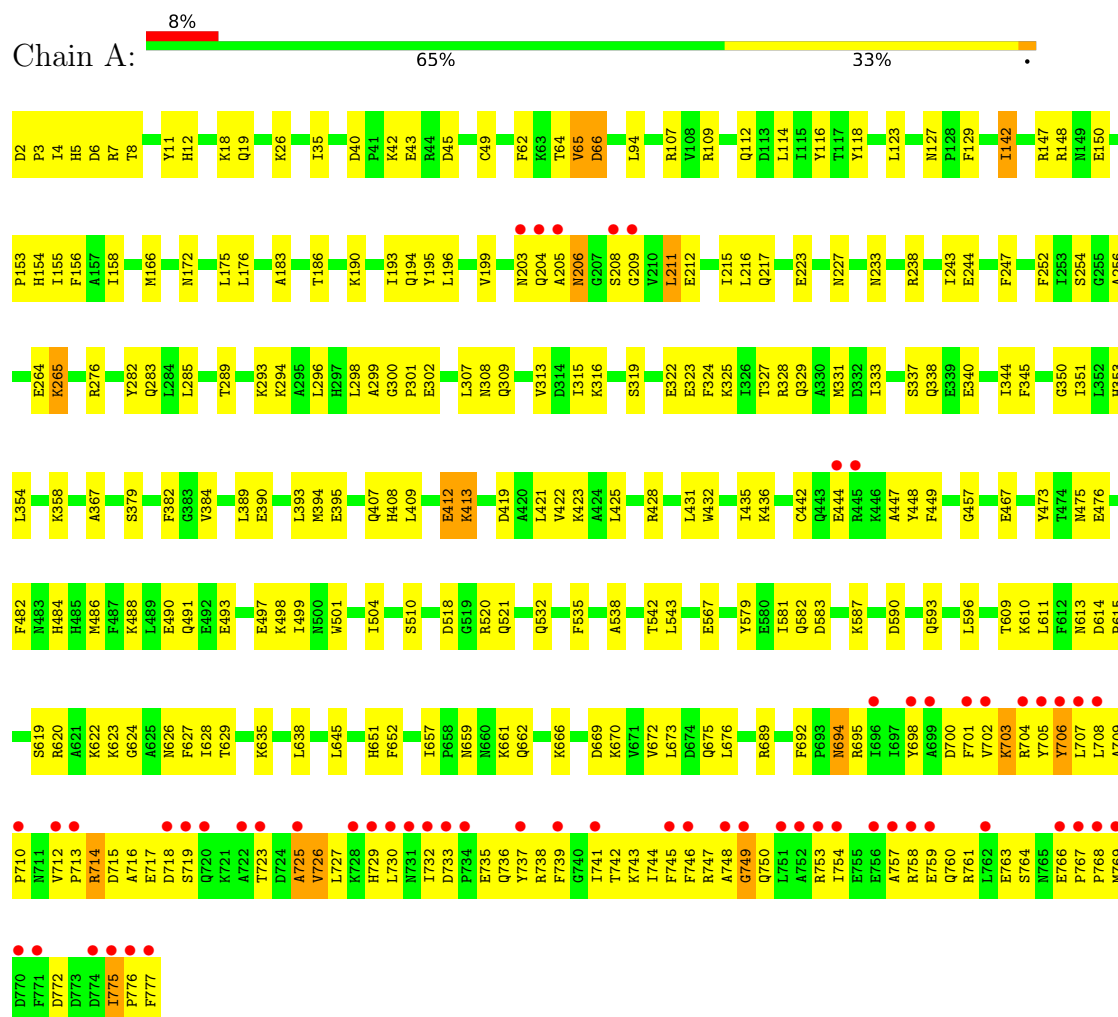
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	435	Total 435	O 435	0	0

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: MYOSIN-2 HEAVY CHAIN



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	88.34Å 146.38Å 152.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.70 – 2.40 19.70 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.8 (19.70-2.40) 99.7 (19.70-2.40)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.07 (at 2.41Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.202 , 0.254 0.203 , 0.208	Depositor DCC
R_{free} test set	1949 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	39.9	Xtriage
Anisotropy	0.357	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 48.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.029 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.036 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6720	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.94% 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: MG, H70, AD9

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	0.57	3/6361 (0.0%)	0.92	20/8587 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	265	LYS	CB-CG	-8.72	1.26	1.52
1	A	265	LYS	CG-CD	-8.71	1.26	1.52
1	A	265	LYS	CD-CE	-6.50	1.32	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	206	ASN	N-CA-C	12.00	131.04	113.40
1	A	703	LYS	N-CA-C	10.58	127.16	112.45
1	A	704	ARG	N-CA-C	-9.93	100.32	114.12
1	A	725	ALA	N-CA-C	9.63	121.55	111.14
1	A	413	LYS	N-CA-C	-8.82	102.09	113.12
1	A	66	ASP	N-CA-C	-7.65	98.00	109.86
1	A	412	GLU	CB-CA-C	-7.54	95.42	110.42
1	A	412	GLU	N-CA-C	6.99	125.69	110.80
1	A	769	MET	CB-CA-C	6.95	124.25	110.42
1	A	703	LYS	CB-CA-C	-6.86	97.15	110.11
1	A	66	ASP	CB-CA-C	-6.49	97.69	113.15
1	A	706	TYR	N-CA-C	-5.96	97.63	108.65
1	A	233	ASN	CB-CA-C	5.93	120.54	109.86
1	A	337	SER	CB-CA-C	-5.92	100.16	109.70
1	A	142	ILE	CB-CA-C	-5.90	104.31	112.04
1	A	447	ALA	N-CA-C	-5.34	106.45	113.12
1	A	717	GLU	N-CA-C	-5.29	107.00	113.50
1	A	300	GLY	CA-C-N	5.20	126.34	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	300	GLY	C-N-CA	5.20	126.34	119.84
1	A	749	GLY	N-CA-C	5.12	125.33	113.18

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	6236	0	6168	322	1
2	A	17	0	4	17	0
3	A	31	0	13	2	0
4	A	1	0	0	0	0
5	A	435	0	0	81	0
All	All	6720	0	6185	327	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 26.

All (327) 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:761:ARG:HD2	1:A:764:SER:CB	1.23	1.61
1:A:761:ARG:CD	1:A:764:SER:HB3	1.33	1.57
1:A:743:LYS:HE3	5:A:2428:HOH:O	1.25	1.34
1:A:761:ARG:CD	1:A:764:SER:CB	1.94	1.28
1:A:700:ASP:HB2	5:A:2414:HOH:O	1.12	1.26
1:A:761:ARG:NH1	1:A:764:SER:HB2	1.54	1.18
1:A:294:LYS:HA	5:A:2186:HOH:O	1.46	1.12
1:A:265:LYS:NZ	2:A:1778:H70:H11	1.47	1.11
1:A:42:LYS:HA	5:A:2038:HOH:O	1.52	1.09
2:A:1778:H70:CL1	5:A:2268:HOH:O	2.10	1.04
1:A:265:LYS:HE3	2:A:1778:H70:HD2	1.06	1.03
1:A:265:LYS:CE	2:A:1778:H70:HD2	1.87	1.02
1:A:761:ARG:CD	1:A:764:SER:HB2	1.82	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:761:ARG:HH11	1:A:764:SER:HB2	1.14	1.01
1:A:772:ASP:HB3	1:A:776:PRO:HG3	1.41	1.01
1:A:761:ARG:CZ	1:A:764:SER:HB2	1.91	1.00
1:A:659:ASN:HD22	1:A:661:LYS:H	1.11	0.99
1:A:18:LYS:H	1:A:112:GLN:HE22	1.10	0.97
1:A:154:HIS:CD2	1:A:156:PHE:H	1.81	0.97
1:A:265:LYS:NZ	2:A:1778:H70:N11	2.13	0.97
1:A:293:LYS:HD2	5:A:2185:HOH:O	1.65	0.96
1:A:615:PRO:O	1:A:619:SER:HB2	1.66	0.96
1:A:694:ASN:HD21	1:A:749:GLY:HA2	1.28	0.95
1:A:265:LYS:HE3	2:A:1778:H70:CD2	1.96	0.95
1:A:118:TYR:CG	1:A:148:ARG:NH2	2.35	0.94
1:A:107:ARG:HD3	5:A:2091:HOH:O	1.65	0.94
1:A:491:GLN:HE21	1:A:501:TRP:HE1	1.06	0.92
1:A:718:ASP:C	5:A:2423:HOH:O	2.11	0.92
1:A:491:GLN:HE22	1:A:504:ILE:H	1.03	0.91
1:A:265:LYS:HZ2	2:A:1778:H70:H11	1.10	0.91
1:A:118:TYR:HB3	1:A:148:ARG:NH2	1.88	0.88
1:A:64:THR:O	1:A:66:ASP:O	1.92	0.88
1:A:431:LEU:CD1	5:A:2271:HOH:O	2.20	0.88
1:A:204:GLN:HG2	5:A:2149:HOH:O	1.72	0.87
1:A:118:TYR:CD1	1:A:148:ARG:NH2	2.42	0.86
1:A:154:HIS:HD2	1:A:156:PHE:H	0.89	0.86
1:A:206:ASN:OD1	1:A:206:ASN:O	1.92	0.86
1:A:520:ARG:HH11	1:A:521:GLN:HE21	1.22	0.86
2:A:1778:H70:BR3	5:A:2171:HOH:O	2.47	0.85
1:A:657:ILE:H	1:A:675:GLN:HE22	1.18	0.85
1:A:761:ARG:NE	1:A:764:SER:CB	2.40	0.85
1:A:154:HIS:HD2	1:A:156:PHE:N	1.74	0.84
1:A:118:TYR:HB3	1:A:148:ARG:HH22	1.43	0.83
1:A:431:LEU:HG	5:A:2271:HOH:O	1.80	0.82
1:A:431:LEU:HD11	5:A:2271:HOH:O	1.79	0.82
1:A:147:ARG:HD3	1:A:147:ARG:H	1.44	0.81
1:A:431:LEU:HD22	2:A:1778:H70:BR2	2.36	0.81
1:A:294:LYS:HD3	5:A:2186:HOH:O	1.79	0.80
1:A:294:LYS:CD	5:A:2186:HOH:O	2.31	0.79
1:A:747:ARG:HD3	5:A:2413:HOH:O	1.82	0.79
1:A:431:LEU:CG	5:A:2271:HOH:O	2.31	0.79
1:A:609:THR:HG23	1:A:613:ASN:HD22	1.48	0.79
1:A:609:THR:HG23	1:A:613:ASN:ND2	1.98	0.79
1:A:761:ARG:NE	1:A:764:SER:HB2	1.98	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:764:SER:HA	1:A:768:PRO:HG2	1.63	0.78
1:A:741:ILE:HG22	1:A:742:THR:HG23	1.64	0.78
1:A:692:PHE:HB3	1:A:745:PHE:HB3	1.66	0.77
1:A:520:ARG:NH1	1:A:521:GLN:HE21	1.81	0.77
1:A:118:TYR:CB	1:A:148:ARG:NH2	2.47	0.77
1:A:746:PHE:C	5:A:2429:HOH:O	2.28	0.76
1:A:761:ARG:CZ	1:A:764:SER:CB	2.65	0.75
1:A:709:ALA:HB1	1:A:710:PRO:HD2	1.69	0.75
1:A:750:GLN:HE22	1:A:753:ARG:HE	1.35	0.74
1:A:26:LYS:HD2	5:A:2022:HOH:O	1.88	0.74
1:A:42:LYS:O	5:A:2038:HOH:O	2.04	0.74
1:A:331:MET:HE1	1:A:345:PHE:CZ	2.22	0.74
1:A:294:LYS:CA	5:A:2186:HOH:O	2.15	0.74
1:A:206:ASN:N	5:A:2153:HOH:O	2.21	0.74
1:A:276:ARG:NH1	5:A:2180:HOH:O	2.22	0.73
1:A:442:CYS:HB3	5:A:2274:HOH:O	1.89	0.73
1:A:294:LYS:CB	5:A:2186:HOH:O	2.34	0.73
1:A:491:GLN:NE2	1:A:504:ILE:H	1.81	0.73
1:A:5:HIS:HD2	5:A:2001:HOH:O	1.71	0.73
1:A:761:ARG:HG3	1:A:763:GLU:O	1.88	0.73
1:A:733:ASP:HB3	1:A:736:GLN:HB2	1.69	0.72
1:A:172:ASN:ND2	1:A:449:PHE:H	1.88	0.71
1:A:694:ASN:ND2	1:A:749:GLY:HA2	2.05	0.71
1:A:491:GLN:HE22	1:A:504:ILE:N	1.85	0.71
1:A:45:ASP:HB3	1:A:673:LEU:HD12	1.73	0.70
1:A:42:LYS:CA	5:A:2038:HOH:O	2.22	0.70
1:A:719:SER:N	5:A:2423:HOH:O	2.21	0.70
1:A:431:LEU:HD21	5:A:2381:HOH:O	1.92	0.69
1:A:421:LEU:O	1:A:425:LEU:HG	1.92	0.69
1:A:610:LYS:HE3	1:A:614:ASP:OD2	1.92	0.69
1:A:659:ASN:ND2	1:A:661:LYS:H	1.87	0.69
1:A:761:ARG:CD	1:A:763:GLU:O	2.41	0.69
1:A:623:LYS:HE2	1:A:628:ILE:HA	1.75	0.68
1:A:743:LYS:NZ	5:A:2427:HOH:O	2.26	0.68
1:A:293:LYS:NZ	5:A:2185:HOH:O	2.26	0.68
1:A:147:ARG:HE	1:A:150:GLU:HB2	1.56	0.68
1:A:5:HIS:CD2	5:A:2001:HOH:O	2.44	0.68
1:A:65:VAL:O	5:A:2060:HOH:O	2.11	0.67
1:A:331:MET:HE1	1:A:345:PHE:HZ	1.56	0.67
1:A:518:ASP:HB2	1:A:635:LYS:HE3	1.77	0.67
1:A:718:ASP:CG	5:A:2423:HOH:O	2.38	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:709:ALA:HA	1:A:729:HIS:CD2	2.30	0.66
1:A:205:ALA:HB1	5:A:2150:HOH:O	1.95	0.66
1:A:316:LYS:HG3	5:A:2197:HOH:O	1.96	0.66
1:A:610:LYS:HD2	1:A:614:ASP:HB2	1.77	0.65
1:A:285:LEU:O	5:A:2181:HOH:O	2.14	0.65
1:A:293:LYS:CD	5:A:2185:HOH:O	2.32	0.65
1:A:206:ASN:CA	5:A:2153:HOH:O	2.44	0.65
1:A:750:GLN:NE2	1:A:753:ARG:HE	1.95	0.65
1:A:761:ARG:NH1	1:A:764:SER:CB	2.47	0.65
1:A:294:LYS:HE2	5:A:2186:HOH:O	1.98	0.64
1:A:325:LYS:O	1:A:329:GLN:HG3	1.98	0.64
1:A:758:ARG:HA	1:A:761:ARG:HH21	1.61	0.63
1:A:26:LYS:CD	5:A:2022:HOH:O	2.44	0.63
1:A:714:ARG:NH2	5:A:2418:HOH:O	2.29	0.63
2:A:1778:H70:C3	5:A:2268:HOH:O	2.47	0.63
1:A:623:LYS:HD3	1:A:628:ILE:HG13	1.80	0.63
1:A:748:ALA:HB3	1:A:753:ARG:NH2	2.13	0.62
2:A:1778:H70:H3	5:A:2268:HOH:O	1.99	0.62
1:A:285:LEU:HD11	1:A:307:LEU:CD1	2.30	0.62
1:A:532:GLN:NE2	1:A:542:THR:C	2.58	0.62
1:A:623:LYS:HZ3	1:A:629:THR:H	1.48	0.62
1:A:294:LYS:CE	5:A:2186:HOH:O	2.48	0.61
1:A:276:ARG:NH2	1:A:282:TYR:CD2	2.67	0.61
1:A:193:ILE:HD11	1:A:243:ILE:HD13	1.81	0.61
1:A:285:LEU:CD1	1:A:307:LEU:CD1	2.79	0.61
1:A:358:LYS:HE3	5:A:2220:HOH:O	2.01	0.61
1:A:730:LEU:H	1:A:758:ARG:HH22	1.48	0.61
1:A:761:ARG:HD3	1:A:764:SER:CB	2.20	0.61
1:A:491:GLN:HE21	1:A:501:TRP:NE1	1.90	0.60
1:A:324:PHE:O	1:A:328:ARG:HG3	2.02	0.60
1:A:195:TYR:CZ	1:A:199:VAL:HG21	2.37	0.59
1:A:775:ILE:N	1:A:776:PRO:CD	2.65	0.59
1:A:431:LEU:HD13	2:A:1778:H70:BR2	2.57	0.59
1:A:761:ARG:CG	1:A:763:GLU:O	2.51	0.59
1:A:18:LYS:H	1:A:112:GLN:NE2	1.91	0.59
1:A:203:ASN:HB2	5:A:2153:HOH:O	2.01	0.58
1:A:254:SER:HB2	1:A:444:GLU:HG2	1.84	0.58
1:A:289:THR:HG21	5:A:2183:HOH:O	2.04	0.58
1:A:520:ARG:NH1	1:A:521:GLN:NE2	2.50	0.58
1:A:45:ASP:CB	1:A:673:LEU:HD12	2.34	0.58
1:A:763:GLU:HA	1:A:767:PRO:HG2	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:657:ILE:N	1:A:675:GLN:HE22	1.96	0.57
1:A:738:ARG:O	1:A:744:ILE:HD12	2.04	0.57
1:A:276:ARG:HH11	1:A:315:ILE:HD11	1.69	0.57
1:A:623:LYS:NZ	1:A:629:THR:H	2.02	0.57
1:A:209:GLY:O	1:A:212:GLU:OE1	2.23	0.57
1:A:428:ARG:NH2	2:A:1778:H70:CL2	2.67	0.57
1:A:730:LEU:HB3	1:A:758:ARG:HH12	1.70	0.57
1:A:175:LEU:HD13	1:A:651:HIS:HB2	1.87	0.56
1:A:276:ARG:NH1	1:A:315:ILE:HD11	2.19	0.56
1:A:293:LYS:CE	5:A:2185:HOH:O	2.53	0.56
1:A:532:GLN:HE21	1:A:543:LEU:N	2.03	0.56
1:A:620:ARG:HG2	1:A:627:PHE:HB3	1.87	0.56
1:A:590:ASP:OD1	2:A:1778:H70:CL1	2.60	0.56
1:A:767:PRO:HB2	1:A:768:PRO:HD3	1.86	0.56
1:A:289:THR:CG2	5:A:2183:HOH:O	2.53	0.56
1:A:730:LEU:H	1:A:758:ARG:NH2	2.03	0.56
1:A:389:LEU:O	1:A:393:LEU:HG	2.05	0.56
1:A:491:GLN:NE2	1:A:501:TRP:HE1	1.89	0.56
1:A:735:GLU:HB2	5:A:2425:HOH:O	2.05	0.56
1:A:172:ASN:HD21	1:A:448:TYR:HA	1.70	0.55
1:A:203:ASN:C	5:A:2153:HOH:O	2.48	0.55
1:A:670:LYS:HB2	5:A:2017:HOH:O	2.07	0.55
1:A:747:ARG:C	5:A:2429:HOH:O	2.50	0.55
1:A:497:GLU:HG3	1:A:745:PHE:HZ	1.71	0.55
1:A:622:LYS:HE3	1:A:624:GLY:O	2.06	0.54
1:A:285:LEU:HD22	1:A:299:ALA:O	2.07	0.54
1:A:26:LYS:HE3	5:A:2022:HOH:O	2.08	0.54
1:A:204:GLN:NE2	5:A:2149:HOH:O	2.32	0.54
1:A:532:GLN:HG3	1:A:538:ALA:HB1	1.90	0.54
1:A:741:ILE:HG22	1:A:742:THR:CG2	2.34	0.54
1:A:94:LEU:O	1:A:689:ARG:NH1	2.41	0.54
1:A:265:LYS:HB3	1:A:423:LYS:HB3	1.89	0.54
1:A:694:ASN:HD22	1:A:746:PHE:HB2	1.71	0.54
1:A:393:LEU:HD23	1:A:596:LEU:HD21	1.89	0.54
1:A:703:LYS:NZ	5:A:2416:HOH:O	2.14	0.54
1:A:707:LEU:HG	1:A:712:VAL:H	1.72	0.53
1:A:209:GLY:HA3	5:A:2145:HOH:O	2.07	0.53
1:A:395:GLU:HA	1:A:407:GLN:O	2.09	0.53
1:A:775:ILE:H	1:A:776:PRO:HD3	1.71	0.53
1:A:623:LYS:N	1:A:626:ASN:O	2.42	0.52
1:A:476:GLU:OE2	1:A:510:SER:HB2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:484:HIS:HD2	1:A:488:LYS:HD3	1.74	0.52
1:A:761:ARG:HD3	1:A:764:SER:HB2	1.81	0.52
1:A:510:SER:HA	5:A:2309:HOH:O	2.09	0.52
1:A:206:ASN:O	1:A:206:ASN:CG	2.53	0.52
1:A:761:ARG:HD2	1:A:763:GLU:O	2.10	0.52
1:A:319:SER:O	1:A:323:GLU:HG2	2.10	0.52
1:A:763:GLU:HG3	1:A:767:PRO:HB2	1.92	0.52
1:A:698:TYR:CD2	1:A:723:THR:HG21	2.45	0.52
1:A:766:GLU:N	1:A:767:PRO:CD	2.72	0.52
1:A:302:GLU:H	1:A:302:GLU:CD	2.18	0.51
1:A:532:GLN:NE2	1:A:542:THR:OG1	2.43	0.51
1:A:265:LYS:CE	2:A:1778:H70:H11	2.19	0.51
1:A:432:TRP:CZ2	1:A:436:LYS:HE3	2.45	0.51
1:A:247:PHE:HA	1:A:252:PHE:O	2.10	0.51
1:A:497:GLU:HG3	1:A:745:PHE:CZ	2.46	0.51
1:A:535:PHE:HD2	1:A:538:ALA:HB2	1.74	0.51
1:A:735:GLU:HG3	5:A:2425:HOH:O	2.11	0.51
1:A:746:PHE:HB3	1:A:750:GLN:HB2	1.93	0.51
1:A:409:LEU:HB3	1:A:413:LYS:HB3	1.93	0.50
1:A:338:GLN:HG3	5:A:2213:HOH:O	2.11	0.50
1:A:436:LYS:HG3	5:A:2272:HOH:O	2.11	0.50
1:A:710:PRO:O	1:A:712:VAL:HG12	2.12	0.50
1:A:739:PHE:CD1	1:A:744:ILE:HD13	2.46	0.50
1:A:467:GLU:OE2	1:A:587:LYS:HE2	2.12	0.50
1:A:2:ASP:OD1	1:A:3:PRO:HD2	2.11	0.50
1:A:223:GLU:O	1:A:227:ASN:HB2	2.12	0.50
1:A:614:ASP:OD1	1:A:614:ASP:C	2.55	0.49
1:A:692:PHE:CB	1:A:745:PHE:HB3	2.38	0.49
1:A:705:TYR:HB3	1:A:726:VAL:HG11	1.94	0.49
1:A:723:THR:O	1:A:727:LEU:HD13	2.11	0.49
1:A:109:ARG:HB3	1:A:114:LEU:HB2	1.93	0.49
1:A:6:ASP:C	1:A:8:THR:H	2.19	0.49
1:A:673:LEU:HD23	1:A:676:LEU:HD12	1.94	0.49
1:A:532:GLN:HE22	1:A:542:THR:C	2.20	0.49
1:A:4:ILE:HA	1:A:11:TYR:CD2	2.48	0.49
1:A:147:ARG:HD3	1:A:147:ARG:N	2.19	0.49
1:A:166:MET:SD	1:A:247:PHE:CD2	3.06	0.49
1:A:205:ALA:CA	5:A:2150:HOH:O	2.61	0.49
1:A:672:VAL:O	1:A:676:LEU:HG	2.12	0.49
1:A:532:GLN:NE2	1:A:543:LEU:N	2.61	0.49
1:A:186:THR:O	1:A:190:LYS:HG3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:LYS:CE	5:A:2022:HOH:O	2.61	0.49
1:A:193:ILE:HD11	1:A:243:ILE:CD1	2.43	0.49
1:A:238:ARG:HD3	1:A:264:GLU:OE1	2.13	0.49
1:A:367:ALA:O	1:A:408:HIS:HE1	1.95	0.48
1:A:457:GLY:N	3:A:1779:AD9:O2G	2.43	0.48
1:A:147:ARG:H	1:A:147:ARG:CD	2.18	0.48
1:A:764:SER:CA	1:A:768:PRO:HG2	2.40	0.48
1:A:211:LEU:O	1:A:215:ILE:HG13	2.12	0.48
1:A:694:ASN:HD21	1:A:749:GLY:CA	2.14	0.48
1:A:7:ARG:NH1	1:A:19:GLN:OE1	2.46	0.48
1:A:350:GLY:HA3	1:A:382:PHE:CZ	2.49	0.48
1:A:40:ASP:OD1	1:A:42:LYS:HB2	2.14	0.48
1:A:354:LEU:HD11	1:A:425:LEU:HD11	1.95	0.48
1:A:718:ASP:CB	5:A:2423:HOH:O	2.60	0.48
1:A:757:ALA:O	1:A:761:ARG:NE	2.47	0.48
1:A:709:ALA:HB2	1:A:726:VAL:HA	1.96	0.48
1:A:43:GLU:CD	5:A:2040:HOH:O	2.58	0.47
1:A:351:ILE:HG23	1:A:422:VAL:HG13	1.96	0.47
1:A:695:ARG:O	1:A:695:ARG:HG2	2.15	0.47
1:A:350:GLY:O	1:A:354:LEU:HG	2.13	0.47
1:A:567:GLU:HA	1:A:579:TYR:O	2.15	0.47
1:A:282:TYR:CE1	1:A:307:LEU:HD22	2.50	0.47
1:A:692:PHE:CZ	1:A:747:ARG:NH2	2.83	0.47
1:A:129:PHE:CE2	1:A:662:GLN:HA	2.50	0.46
1:A:733:ASP:OD1	1:A:735:GLU:HG2	2.16	0.46
1:A:276:ARG:NH2	1:A:282:TYR:CG	2.83	0.46
1:A:217:GLN:HG3	1:A:333:ILE:HG21	1.97	0.46
1:A:296:LEU:HB2	1:A:298:LEU:HG	1.98	0.46
1:A:749:GLY:N	5:A:2429:HOH:O	2.47	0.46
1:A:775:ILE:N	1:A:776:PRO:HD3	2.29	0.46
1:A:244:GLU:O	1:A:256:ALA:HA	2.16	0.46
1:A:204:GLN:HG2	5:A:2147:HOH:O	1.77	0.46
1:A:390:GLU:HG3	1:A:394:MET:CE	2.45	0.46
1:A:457:GLY:HA2	1:A:475:ASN:HD21	1.80	0.46
1:A:709:ALA:CB	1:A:726:VAL:HA	2.46	0.46
1:A:296:LEU:O	1:A:353:HIS:HE1	1.99	0.46
1:A:490:GLU:O	1:A:493:GLU:HB3	2.15	0.46
1:A:535:PHE:CD2	1:A:538:ALA:HB2	2.51	0.46
1:A:777:PHE:CD1	1:A:777:PHE:OXT	2.69	0.46
1:A:26:LYS:NZ	5:A:2024:HOH:O	2.47	0.46
1:A:498:LYS:HD3	1:A:741:ILE:HD11	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:ASP:OD1	1:A:8:THR:HB	2.14	0.46
1:A:254:SER:HB2	1:A:444:GLU:CG	2.46	0.46
1:A:328:ARG:NH2	5:A:2208:HOH:O	2.49	0.46
1:A:340:GLU:O	1:A:344:ILE:HG13	2.16	0.45
1:A:205:ALA:CB	5:A:2150:HOH:O	2.57	0.45
1:A:761:ARG:CD	1:A:763:GLU:C	2.89	0.45
1:A:766:GLU:H	1:A:767:PRO:CD	2.28	0.45
1:A:301:PRO:HB2	1:A:313:VAL:HG11	1.98	0.45
1:A:666:LYS:NZ	5:A:2400:HOH:O	2.49	0.45
1:A:484:HIS:CD2	1:A:488:LYS:HD3	2.52	0.45
1:A:712:VAL:O	1:A:712:VAL:HG13	2.17	0.45
1:A:116:TYR:HB3	1:A:123:LEU:HD11	1.99	0.44
1:A:155:ILE:HD12	1:A:158:ILE:CG2	2.47	0.44
1:A:726:VAL:HG13	1:A:727:LEU:HD12	1.98	0.44
1:A:761:ARG:HD3	1:A:763:GLU:C	2.43	0.44
1:A:206:ASN:HA	5:A:2153:HOH:O	2.15	0.44
1:A:354:LEU:CD1	1:A:425:LEU:CD1	2.95	0.44
1:A:581:ILE:O	1:A:582:GLN:C	2.58	0.44
1:A:743:LYS:CE	5:A:2428:HOH:O	2.10	0.44
1:A:706:TYR:O	1:A:708:LEU:N	2.48	0.44
1:A:763:GLU:O	1:A:764:SER:HB3	2.18	0.44
1:A:419:ASP:O	1:A:423:LYS:HG3	2.18	0.44
1:A:118:TYR:CE2	1:A:153:PRO:HA	2.53	0.43
1:A:194:GLN:HG3	5:A:2141:HOH:O	2.17	0.43
1:A:283:GLN:NE2	1:A:327:THR:OG1	2.46	0.43
1:A:308:ASN:OD1	1:A:309:GLN:NE2	2.50	0.43
1:A:344:ILE:HD11	1:A:432:TRP:HZ3	1.83	0.43
1:A:457:GLY:H	3:A:1779:AD9:H2G1	1.64	0.43
1:A:759:GLU:HB3	1:A:760:GLN:OE1	2.18	0.43
1:A:49:CYS:SG	5:A:2051:HOH:O	2.62	0.43
1:A:701:PHE:HZ	1:A:727:LEU:HD11	1.83	0.43
1:A:744:ILE:O	1:A:744:ILE:HG23	2.19	0.43
1:A:282:TYR:CZ	1:A:307:LEU:HD22	2.54	0.43
1:A:750:GLN:HE22	1:A:753:ARG:NE	2.11	0.43
1:A:709:ALA:HA	1:A:729:HIS:HD2	1.80	0.42
1:A:35:ILE:HG21	1:A:62:PHE:CG	2.54	0.42
1:A:127:ASN:HB2	1:A:183:ALA:O	2.19	0.42
1:A:176:LEU:HD12	1:A:652:PHE:CE1	2.54	0.42
1:A:206:ASN:HB2	5:A:2153:HOH:O	2.19	0.42
1:A:285:LEU:HD11	1:A:307:LEU:HD11	1.98	0.42
1:A:354:LEU:CD1	1:A:425:LEU:HD11	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:473:TYR:CE2	1:A:638:LEU:HD23	2.53	0.42
1:A:736:GLN:OE1	1:A:747:ARG:HD2	2.19	0.42
1:A:767:PRO:CB	1:A:768:PRO:HD3	2.49	0.42
1:A:582:GLN:O	1:A:583:ASP:HB2	2.20	0.42
1:A:6:ASP:C	1:A:8:THR:N	2.78	0.42
1:A:593:GLN:NE2	5:A:2373:HOH:O	2.52	0.42
1:A:723:THR:C	1:A:725:ALA:H	2.28	0.42
1:A:745:PHE:N	1:A:745:PHE:CD2	2.88	0.42
1:A:379:SER:HB3	1:A:384:VAL:O	2.20	0.41
1:A:669:ASP:O	1:A:673:LEU:HG	2.20	0.41
1:A:727:LEU:HD23	1:A:737:TYR:CE2	2.55	0.41
1:A:294:LYS:HB2	5:A:2186:HOH:O	2.09	0.41
1:A:482:PHE:CZ	1:A:486:MET:HG3	2.55	0.41
1:A:610:LYS:CD	1:A:614:ASP:HB2	2.48	0.41
1:A:206:ASN:CB	5:A:2153:HOH:O	2.68	0.41
1:A:763:GLU:HG3	1:A:764:SER:H	1.86	0.41
1:A:45:ASP:CG	1:A:673:LEU:HD12	2.46	0.41
1:A:422:VAL:HA	1:A:425:LEU:HD12	2.03	0.41
1:A:707:LEU:HD11	1:A:713:PRO:HG3	2.03	0.41
1:A:732:ILE:HG22	1:A:754:ILE:HG12	2.02	0.41
1:A:737:TYR:HE1	1:A:739:PHE:CE2	2.39	0.41
2:A:1778:H70:CE2	5:A:2268:HOH:O	2.53	0.41
1:A:622:LYS:HE2	1:A:622:LYS:HB3	1.88	0.41
1:A:3:PRO:HB2	1:A:142:ILE:CD1	2.51	0.40
1:A:520:ARG:HH11	1:A:520:ARG:HG2	1.87	0.40
1:A:715:ASP:O	1:A:716:ALA:HB2	2.21	0.40
1:A:431:LEU:HB2	2:A:1778:H70:BR2	2.77	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:A:204:GLN:OE1	1:A:615:PRO:CB[3_754]	1.76	0.44

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	774/776 (100%)	717 (93%)	52 (7%)	5 (1%)	21	32

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	714	ARG
1	A	208	SER
1	A	702	VAL
1	A	775	ILE
1	A	499	ILE

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	680/680 (100%)	668 (98%)	12 (2%)	51	73

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	HIS
1	A	65	VAL
1	A	196	LEU
1	A	211	LEU
1	A	216	LEU
1	A	322	GLU
1	A	412	GLU
1	A	435	ILE
1	A	611	LEU
1	A	645	LEU
1	A	694	ASN
1	A	726	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
1	A	104	HIS
1	A	112	GLN
1	A	154	HIS
1	A	172	ASN
1	A	194	GLN
1	A	227	ASN
1	A	283	GLN
1	A	353	HIS
1	A	408	HIS
1	A	475	ASN
1	A	484	HIS
1	A	491	GLN
1	A	511	GLN
1	A	521	GLN
1	A	532	GLN
1	A	541	ASN
1	A	593	GLN
1	A	613	ASN
1	A	649	ASN
1	A	659	ASN
1	A	675	GLN
1	A	679	ASN
1	A	694	ASN
1	A	729	HIS
1	A	750	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

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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
3	AD9	A	1779	4	29,33,33	1.94	6 (20%)	42,52,52	1.79	11 (26%)
2	H70	A	1778	-	17,18,18	2.01	6 (35%)	22,27,27	1.13	1 (4%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	AD9	A	1779	4	-	0/16/38/38	0/3/3/3
2	H70	A	1778	-	-	3/4/4/4	0/2/2/2

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1779	AD9	O3B-VG	6.66	2.10	1.80
2	A	1778	H70	CR-CH	4.08	1.45	1.39
3	A	1779	AD9	C2'-C3'	-4.02	1.42	1.53
3	A	1779	AD9	C6-N6	3.72	1.43	1.34
2	A	1778	H70	C3-CE2	3.58	1.44	1.38
2	A	1778	H70	C3-CH	3.17	1.43	1.38
3	A	1779	AD9	C3'-C4'	-2.52	1.46	1.53
3	A	1779	AD9	O5'-C5'	-2.26	1.36	1.44
2	A	1778	H70	CG-CR	2.20	1.45	1.41
2	A	1778	H70	C9-C8	2.12	1.45	1.38
3	A	1779	AD9	C5-N7	-2.11	1.35	1.39
2	A	1778	H70	BR2-C8	-2.11	1.83	1.88

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1779	AD9	N3-C2-N1	-5.32	120.53	128.58
3	A	1779	AD9	C5-C4-N3	-3.85	121.41	126.72
3	A	1779	AD9	C2-N1-C6	3.27	124.11	118.73
3	A	1779	AD9	N9-C8-N7	-3.10	109.53	113.94
3	A	1779	AD9	C2-N3-C4	3.06	119.30	111.83
3	A	1779	AD9	O4'-C1'-N9	2.88	113.62	108.09
3	A	1779	AD9	C4-N9-C8	2.73	108.60	105.74
3	A	1779	AD9	N3-C4-N9	2.72	131.79	127.17
3	A	1779	AD9	O2B-PB-O3A	2.68	114.52	107.27
3	A	1779	AD9	C5-N7-C8	2.37	107.18	103.45
2	A	1778	H70	CR-CH-CL2	2.35	121.73	118.81
3	A	1779	AD9	C4-C5-N7	-2.34	107.91	110.58

There are no chirality outliers.

All (3) torsion outliers are listed below:

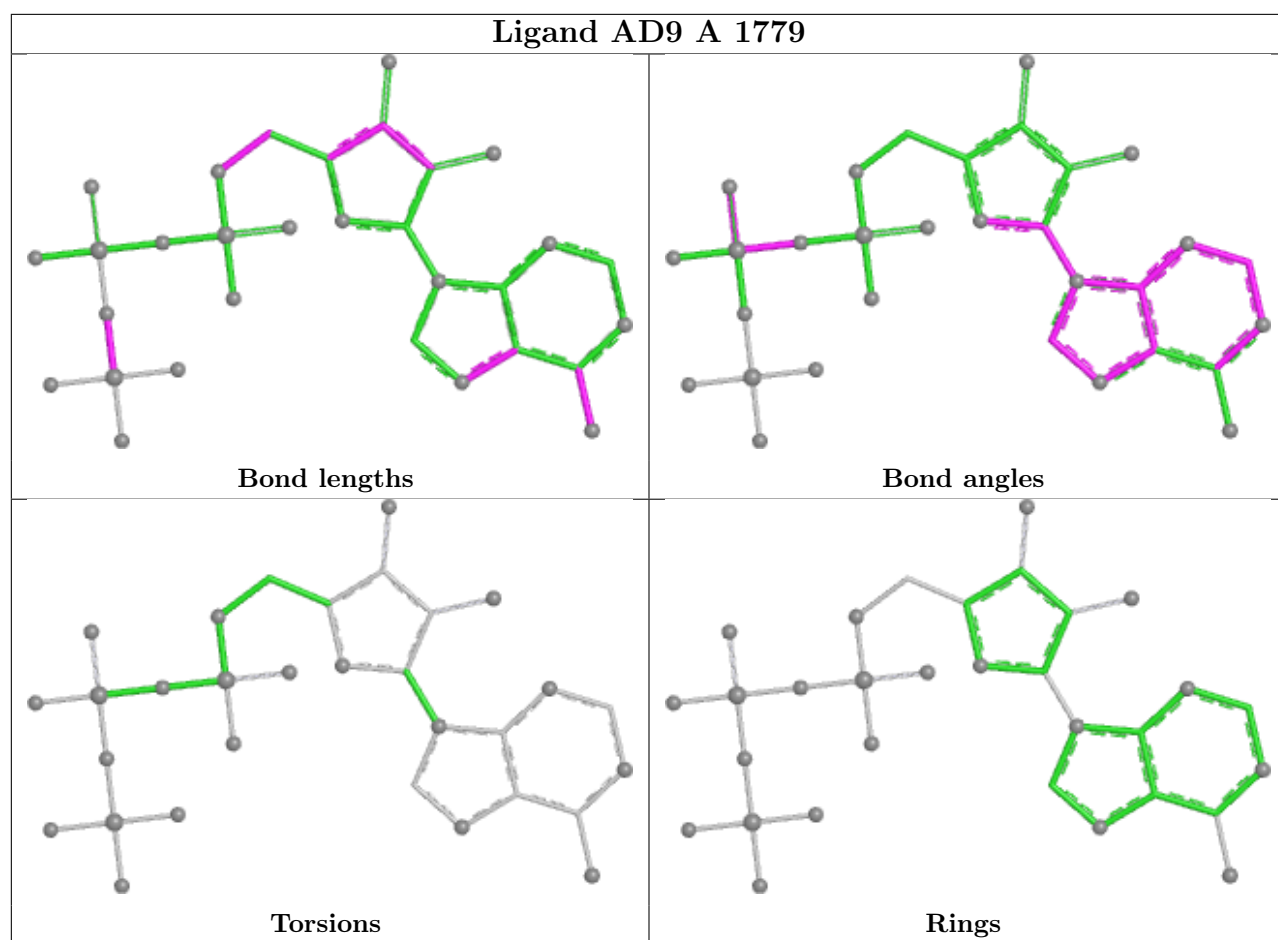
Mol	Chain	Res	Type	Atoms
2	A	1778	H70	C9-CD1-CG-CD2
2	A	1778	H70	N11-CD1-CG-CD2
2	A	1778	H70	C9-CD1-CG-CR

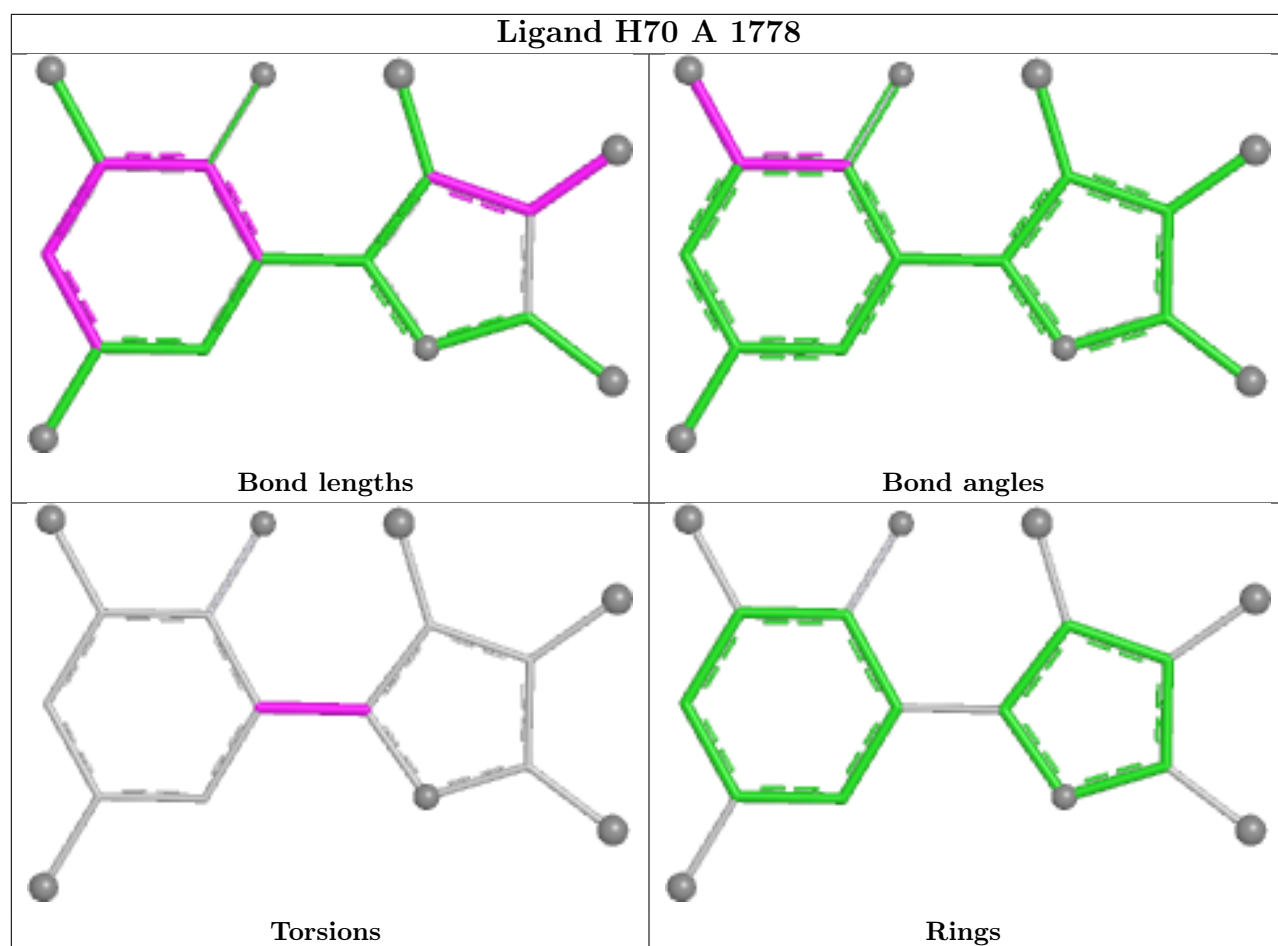
There are no ring outliers.

2 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1779	AD9	2	0
2	A	1778	H70	17	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.





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	776/776 (100%)	0.07	59 (7%)	20 16	23, 43, 119, 157	21 (2%)

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	757	ALA	6.0
1	A	762	LEU	5.8
1	A	777	PHE	5.5
1	A	702	VAL	4.9
1	A	754	ILE	4.6
1	A	730	LEU	4.6
1	A	771	PHE	4.5
1	A	710	PRO	4.3
1	A	776	PRO	4.1
1	A	775	ILE	4.0
1	A	706	TYR	4.0
1	A	756	GLU	3.9
1	A	204	GLN	3.8
1	A	768	PRO	3.7
1	A	712	VAL	3.6
1	A	752	ALA	3.4
1	A	705	TYR	3.3
1	A	707	LEU	3.3
1	A	769	MET	3.3
1	A	737	TYR	3.2
1	A	770	ASP	3.2
1	A	719	SER	3.2
1	A	698	TYR	3.2
1	A	720	GLN	3.0
1	A	723	THR	3.0
1	A	729	HIS	2.9
1	A	758	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	774	ASP	2.9
1	A	696	ILE	2.8
1	A	699	ALA	2.7
1	A	208	SER	2.7
1	A	701	PHE	2.6
1	A	753	ARG	2.6
1	A	722	ALA	2.6
1	A	728	LYS	2.6
1	A	734	PRO	2.6
1	A	751	LEU	2.6
1	A	767	PRO	2.6
1	A	741	ILE	2.5
1	A	444	GLU	2.4
1	A	759	GLU	2.4
1	A	445	ARG	2.4
1	A	766	GLU	2.3
1	A	203	ASN	2.3
1	A	718	ASP	2.3
1	A	205	ALA	2.2
1	A	725	ALA	2.2
1	A	732	ILE	2.2
1	A	745	PHE	2.2
1	A	739	PHE	2.2
1	A	731	ASN	2.1
1	A	748	ALA	2.1
1	A	708	LEU	2.1
1	A	749	GLY	2.1
1	A	746	PHE	2.1
1	A	209	GLY	2.1
1	A	713	PRO	2.1
1	A	704	ARG	2.0
1	A	733	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

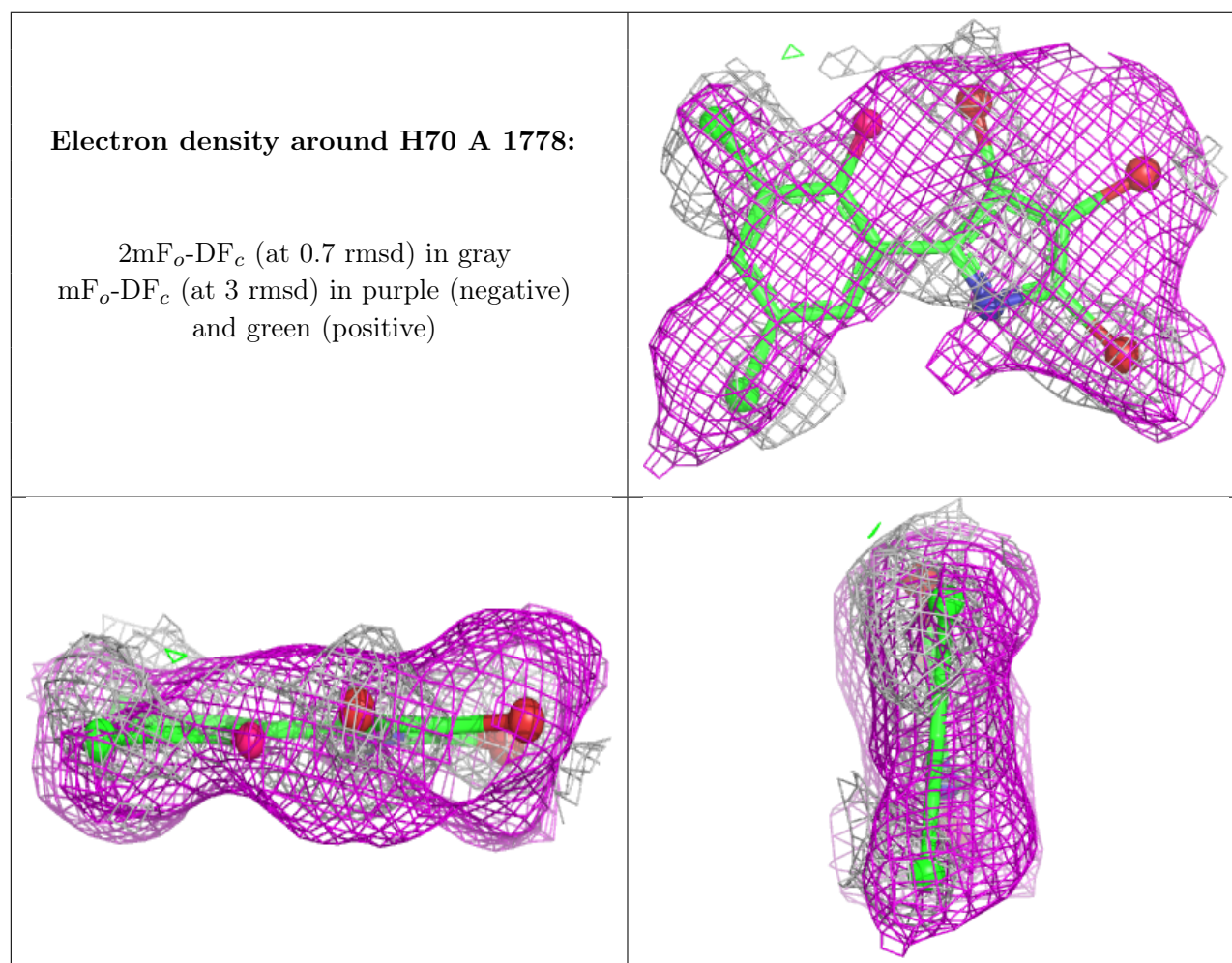
There are no oligosaccharides in this entry.

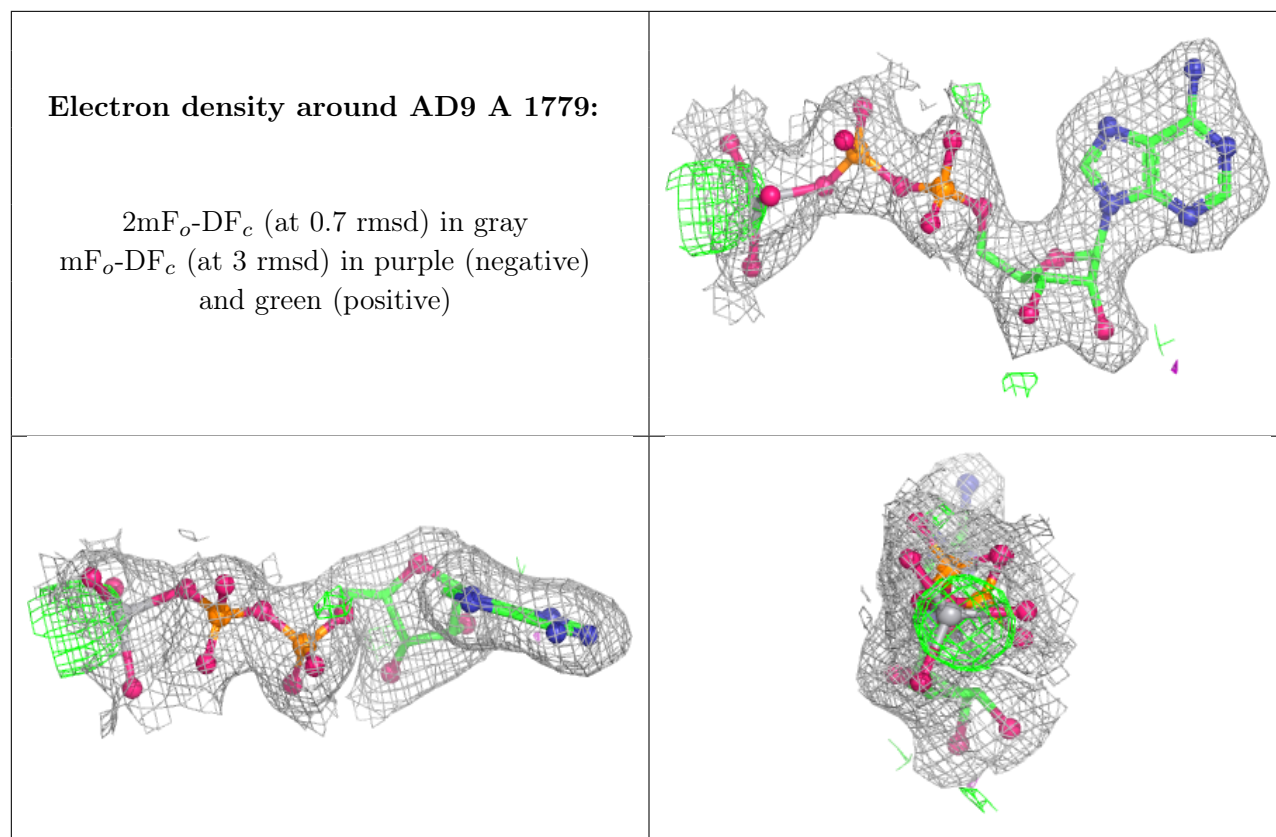
6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	H70	A	1778	17/17	0.66	0.15	56,64,117,123	0
4	MG	A	1780	1/1	0.97	0.07	37,37,37,37	0
3	AD9	A	1779	31/31	0.98	0.06	24,29,33,34	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.