



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:38 AM UTC

PDB ID : 3EQ8 / pdb_00003eq8
Title : Prolyl oligopeptidase complexed with R-Pro-(decarboxy-Pro)-Type inhibitors
Authors : Kanai, K.; Aranyi, P.; Bocskei, Z.; Ferenczy, G.; Harmat, V.; Simon, K.;
Naray-Szabo, G.; Hermecz, I.
Deposited on : 2008-09-30
Resolution : 2.73 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

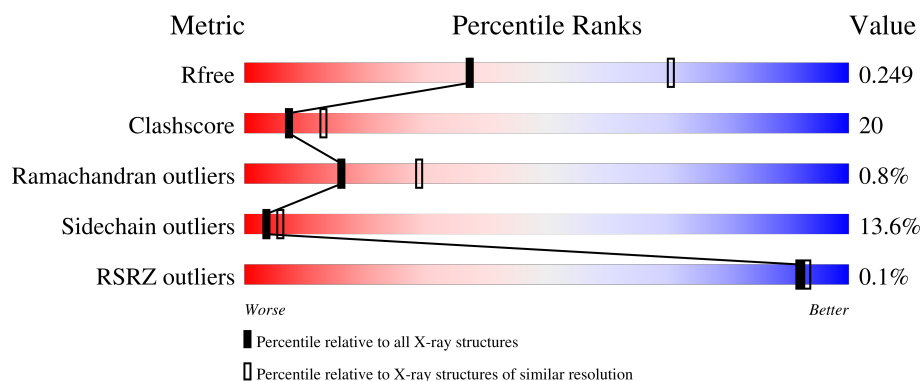
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.73 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	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
Sidechain outliers	187428	1831 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

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	710	60% 33% 6%

2 Entry composition [i](#)

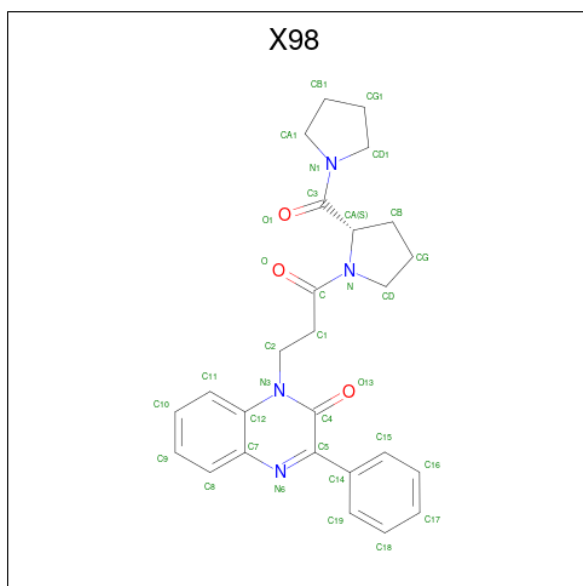
There are 3 unique types of molecules in this entry. The entry contains 5788 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Prolyl endopeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	709	Total	C	N	O	S	0	3	0
			5656	3631	936	1062	27			

- Molecule 2 is 1-{3-oxo-3-[(2S)-2-(pyrrolidin-1-ylcarbonyl)pyrrolidin-1-yl]propyl}-3-phenylquinoxalin-2(1H)-one (CCD ID: X98) (formula: C₂₆H₂₈N₄O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			33	26	4	3		

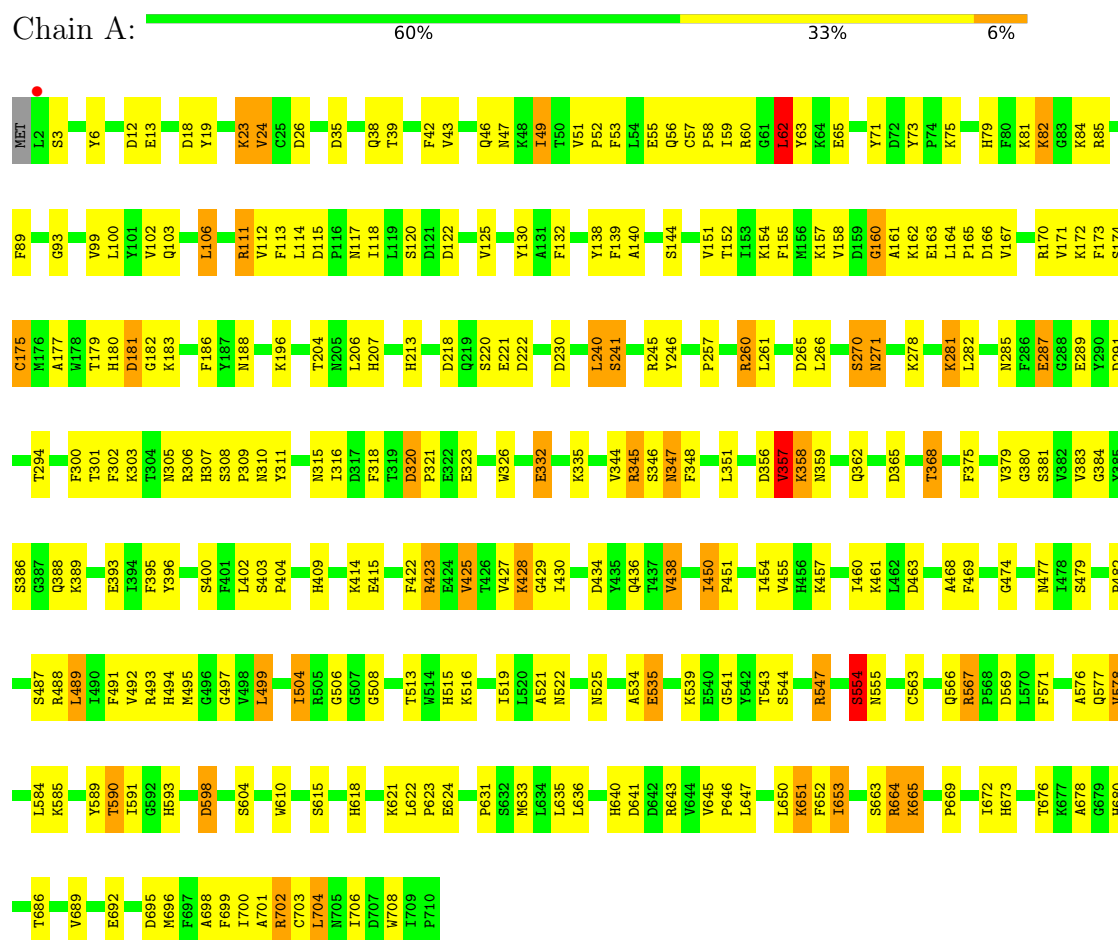
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	99	Total	0	0
			99		

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: Prolyl endopeptidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.23Å 101.50Å 112.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.03 – 2.73 34.03 – 2.73	Depositor EDS
% Data completeness (in resolution range)	(Not available) (34.03-2.73) 92.8 (34.03-2.73)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 2.72Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.177 , 0.253 0.172 , 0.249	Depositor DCC
R_{free} test set	2108 reflections (10.01%)	wwPDB-VP
Wilson B-factor (Å ²)	39.4	Xtriage
Anisotropy	0.090	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 50.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5788	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% 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: X98

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	0.42	0/5827	0.90	21/7908 (0.3%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	181	ASP	N-CA-C	-8.51	103.06	113.19
1	A	320	ASP	CA-C-N	8.09	128.18	119.28
1	A	320	ASP	C-N-CA	8.09	128.18	119.28
1	A	698	ALA	N-CA-C	-7.81	102.36	111.03
1	A	62	LEU	N-CA-C	-6.44	103.95	110.97
1	A	554	SER	CB-CA-C	-5.75	109.96	116.63
1	A	158	VAL	N-CA-C	5.51	116.90	110.62
1	A	260	ARG	N-CA-C	-5.51	102.36	110.52
1	A	375	PHE	CA-C-N	5.45	125.71	119.83
1	A	375	PHE	C-N-CA	5.45	125.71	119.83
1	A	506	GLY	N-CA-C	-5.43	107.53	114.37
1	A	429	GLY	N-CA-C	-5.42	108.03	114.48
1	A	6	TYR	CA-C-N	5.26	126.00	120.11
1	A	6	TYR	C-N-CA	5.26	126.00	120.11
1	A	160	GLY	N-CA-C	-5.21	107.92	115.27
1	A	26	ASP	CA-C-N	5.20	124.82	119.56
1	A	26	ASP	C-N-CA	5.20	124.82	119.56
1	A	179	THR	N-CA-C	-5.16	102.84	110.48
1	A	175	CYS	N-CA-C	-5.07	102.57	110.42
1	A	615	SER	CA-C-N	5.05	124.49	119.24
1	A	615	SER	C-N-CA	5.05	124.49	119.24

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	5656	0	5417	226	0
2	A	33	0	28	10	0
3	A	99	0	0	8	0
All	All	5788	0	5445	228	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 20.

All (228) 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:60:ARG:HG2	1:A:60:ARG:HH11	1.22	1.02
1:A:428:LYS:H	1:A:428:LYS:HD3	1.22	1.02
1:A:423:ARG:HH11	1:A:423:ARG:CG	1.85	0.88
1:A:423:ARG:HH11	1:A:423:ARG:CB	1.86	0.88
1:A:51:VAL:HB	1:A:52:PRO:HD3	1.56	0.86
1:A:428:LYS:H	1:A:428:LYS:CD	1.89	0.85
1:A:423:ARG:CB	1:A:423:ARG:NH1	2.44	0.80
1:A:547:ARG:HA	1:A:704:LEU:HD13	1.63	0.80
1:A:578:VAL:HG23	1:A:578:VAL:O	1.82	0.78
1:A:257:PRO:HG3	1:A:477:ASN:HD22	1.50	0.77
1:A:696:MET:O	1:A:700:ILE:HG13	1.84	0.77
1:A:59:ILE:HD11	1:A:708:TRP:HB2	1.66	0.77
1:A:647:LEU:HD12	1:A:647:LEU:C	2.11	0.75
1:A:57:CYS:SG	1:A:59:ILE:HG22	2.26	0.75
1:A:423:ARG:HH11	1:A:423:ARG:HG2	1.49	0.74
1:A:438:VAL:HG22	1:A:454:ILE:HB	1.71	0.73
1:A:423:ARG:NH1	1:A:423:ARG:HB2	2.05	0.71
1:A:460:ILE:HG13	1:A:460:ILE:O	1.91	0.71
1:A:646:PRO:O	1:A:650:LEU:HG	1.91	0.70
1:A:60:ARG:HG2	1:A:60:ARG:NH1	1.97	0.70
1:A:436:GLN:HE21	1:A:438:VAL:CG1	2.06	0.69
1:A:241:SER:OG	1:A:246:TYR:HB2	1.93	0.69
1:A:577:GLN:O	1:A:578:VAL:C	2.36	0.68
1:A:344:VAL:O	1:A:348:PHE:HB2	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:566:GLN:O	1:A:567:ARG:HG2	1.94	0.68
1:A:499:LEU:C	1:A:499:LEU:HD23	2.19	0.68
1:A:635:LEU:HB2	1:A:672:ILE:HG13	1.76	0.67
1:A:543:THR:HG23	1:A:544:SER:N	2.09	0.67
1:A:554:SER:OG	2:A:711:X98:HDA	1.95	0.67
1:A:488:ARG:O	1:A:491:PHE:HB3	1.96	0.66
1:A:404:PRO:HB2	1:A:425:VAL:HG22	1.78	0.66
1:A:491:PHE:O	1:A:495:MET:HB2	1.96	0.65
1:A:428:LYS:HD3	1:A:428:LYS:N	2.02	0.64
1:A:47:ASN:O	1:A:51:VAL:HG23	1.98	0.64
1:A:100:LEU:HD23	1:A:113:PHE:HD2	1.63	0.64
1:A:358:LYS:HG2	1:A:379:VAL:HG13	1.80	0.62
1:A:345:ARG:O	1:A:347:ASN:N	2.31	0.62
1:A:305:ASN:HB3	1:A:311:TYR:CE2	2.35	0.62
1:A:120:SER:HB3	1:A:125:VAL:HG23	1.81	0.62
1:A:589:TYR:CE1	1:A:647:LEU:HD21	2.35	0.61
1:A:291:ASP:HB3	1:A:303:LYS:HB3	1.80	0.61
1:A:365:ASP:OD1	1:A:368:THR:HG23	2.00	0.61
1:A:60:ARG:HH11	1:A:60:ARG:CG	2.04	0.61
1:A:664:ARG:HG3	1:A:664:ARG:HH11	1.65	0.61
1:A:59:ILE:CD1	1:A:708:TRP:HB2	2.31	0.61
1:A:323:GLU:HA	1:A:326:TRP:CE2	2.35	0.61
1:A:63:TYR:OH	1:A:487:SER:O	2.18	0.60
1:A:261:LEU:HD12	1:A:282:LEU:HD23	1.83	0.60
1:A:285:ASN:OD1	1:A:287:GLU:HB2	2.00	0.60
1:A:173:PHE:HD2	2:A:711:X98:H19	1.66	0.60
1:A:455:VAL:CG2	1:A:499:LEU:HD22	2.31	0.60
1:A:567:ARG:NH2	1:A:569:ASP:OD2	2.33	0.60
1:A:52:PRO:O	1:A:56:GLN:HG3	2.00	0.60
1:A:563:CYS:O	1:A:571:PHE:HE1	1.85	0.60
1:A:643:ARG:HG2	1:A:680:HIS:CD2	2.37	0.60
1:A:270:SER:C	1:A:271:ASN:OD1	2.45	0.59
1:A:51:VAL:O	1:A:55:GLU:HG3	2.02	0.58
1:A:356:ASP:C	1:A:357:VAL:HG23	2.28	0.58
1:A:635:LEU:O	1:A:636:LEU:HD23	2.04	0.58
1:A:51:VAL:HB	1:A:52:PRO:CD	2.30	0.58
1:A:631:PRO:O	1:A:633:MET:HG3	2.03	0.57
1:A:468:ALA:HB2	1:A:543:THR:HB	1.87	0.56
1:A:555:ASN:HB2	2:A:711:X98:HB	1.86	0.56
1:A:635:LEU:HD11	1:A:652:PHE:HD2	1.69	0.56
1:A:345:ARG:NH2	1:A:348:PHE:CZ	2.73	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:621:LYS:NZ	1:A:621:LYS:HB2	2.19	0.56
1:A:60:ARG:NH1	1:A:60:ARG:CG	2.65	0.56
1:A:423:ARG:CG	1:A:423:ARG:NH1	2.56	0.56
1:A:102:VAL:HG12	1:A:111:ARG:O	2.06	0.55
1:A:347:ASN:N	1:A:347:ASN:HD22	2.05	0.55
1:A:522:ASN:HB3	1:A:525:ASN:ND2	2.22	0.55
1:A:173:PHE:CE2	2:A:711:X98:H2	2.42	0.55
1:A:62:LEU:HD12	1:A:708:TRP:CG	2.42	0.55
1:A:106:LEU:HD11	1:A:422:PHE:HZ	1.72	0.54
1:A:383:VAL:CG1	1:A:383:VAL:O	2.55	0.54
1:A:89:PHE:CE2	1:A:100:LEU:HB2	2.43	0.54
1:A:469:PHE:HB2	1:A:491:PHE:CE2	2.43	0.53
1:A:59:ILE:HG21	1:A:701:ALA:HB3	1.91	0.53
1:A:423:ARG:HB2	1:A:423:ARG:CZ	2.37	0.53
2:A:711:X98:H19	2:A:711:X98:O13	2.09	0.53
1:A:427:VAL:HG12	1:A:427:VAL:O	2.08	0.53
1:A:245:ARG:HG2	1:A:246:TYR:CD1	2.44	0.52
1:A:180:HIS:C	1:A:182:GLY:H	2.17	0.52
1:A:488:ARG:HD2	1:A:499:LEU:HD12	1.91	0.52
1:A:643:ARG:NH1	2:A:711:X98:H1A	2.25	0.52
1:A:59:ILE:HG21	1:A:701:ALA:CB	2.40	0.52
1:A:436:GLN:HE21	1:A:438:VAL:HG12	1.75	0.52
1:A:584:LEU:HD12	1:A:618:HIS:NE2	2.24	0.52
1:A:673:HIS:HD2	3:A:789:HOH:O	1.91	0.52
1:A:164:LEU:HB3	1:A:165:PRO:HD2	1.91	0.51
1:A:388:GLN:HB3	3:A:739:HOH:O	2.10	0.51
1:A:704:LEU:N	1:A:704:LEU:HD23	2.25	0.51
1:A:519:ILE:HG22	1:A:598:ASP:OD2	2.11	0.51
1:A:19:TYR:HD1	1:A:24:VAL:HG13	1.76	0.51
1:A:347:ASN:ND2	1:A:347:ASN:H	2.08	0.51
1:A:593:HIS:HD2	3:A:721:HOH:O	1.93	0.51
1:A:647:LEU:O	1:A:651:LYS:HG3	2.11	0.51
1:A:99:VAL:HG11	1:A:112:VAL:HG13	1.91	0.51
1:A:566:GLN:C	1:A:567:ARG:HG2	2.36	0.51
1:A:18:ASP:OD2	1:A:23:LYS:NZ	2.45	0.50
1:A:174:SER:OG	1:A:175:CYS:N	2.43	0.50
1:A:555:ASN:ND2	3:A:779:HOH:O	2.44	0.50
1:A:75:LYS:HE3	1:A:93:GLY:O	2.12	0.50
1:A:347:ASN:N	1:A:347:ASN:ND2	2.60	0.50
1:A:300:PHE:CE1	1:A:318:PHE:CE1	3.00	0.50
1:A:430:ILE:HD13	1:A:489:LEU:HD12	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:PRO:HB3	1:A:513:THR:HG23	1.94	0.49
1:A:640:HIS:HB2	1:A:678:ALA:O	2.12	0.49
1:A:673:HIS:CD2	3:A:789:HOH:O	2.65	0.49
1:A:539:LYS:C	1:A:541:GLY:H	2.20	0.49
1:A:664:ARG:HG3	1:A:664:ARG:NH1	2.26	0.49
1:A:703:CYS:C	1:A:704:LEU:HD23	2.37	0.49
1:A:173:PHE:CD2	2:A:711:X98:O13	2.66	0.49
1:A:38:GLN:OE1	1:A:38:GLN:N	2.40	0.49
1:A:106:LEU:HD11	1:A:422:PHE:CZ	2.47	0.49
1:A:645:VAL:HG22	3:A:726:HOH:O	2.12	0.49
1:A:653:ILE:HG12	1:A:672:ILE:HB	1.94	0.49
1:A:271:ASN:OD1	1:A:271:ASN:N	2.45	0.49
1:A:493:ARG:NH1	1:A:494:HIS:ND1	2.60	0.49
1:A:636:LEU:HD22	1:A:673:HIS:HB3	1.95	0.49
1:A:477:ASN:OD1	1:A:508:GLY:HA3	2.11	0.49
1:A:240:LEU:CD2	1:A:240:LEU:N	2.75	0.49
1:A:591:ILE:O	1:A:591:ILE:HG13	2.12	0.49
1:A:578:VAL:O	1:A:578:VAL:CG2	2.54	0.48
1:A:306:ARG:O	1:A:307:HIS:HB2	2.12	0.48
1:A:404:PRO:HB2	1:A:425:VAL:CG2	2.43	0.48
1:A:106:LEU:HD12	1:A:106:LEU:O	2.14	0.48
1:A:347:ASN:HD22	1:A:347:ASN:H	1.61	0.48
1:A:699:PHE:O	1:A:703:CYS:HB2	2.14	0.48
1:A:257:PRO:C	1:A:289:GLU:HB2	2.38	0.48
1:A:650:LEU:O	1:A:653:ILE:HG22	2.14	0.48
1:A:310:ASN:HD22	1:A:310:ASN:N	2.12	0.48
1:A:12:ASP:C	1:A:12:ASP:OD1	2.57	0.48
1:A:177:ALA:HB3	1:A:186:PHE:HB2	1.96	0.48
1:A:85:ARG:HG3	3:A:733:HOH:O	2.14	0.47
1:A:174:SER:OG	1:A:175:CYS:O	2.31	0.47
1:A:180:HIS:C	1:A:182:GLY:N	2.69	0.47
1:A:455:VAL:HG23	1:A:499:LEU:HD22	1.95	0.47
1:A:19:TYR:CD1	1:A:24:VAL:HG13	2.49	0.47
1:A:52:PRO:O	1:A:53:PHE:C	2.57	0.47
1:A:316:ILE:HG12	1:A:326:TRP:NE1	2.29	0.47
1:A:457:LYS:O	1:A:460:ILE:HG23	2.15	0.47
1:A:516:LYS:HA	1:A:519:ILE:HG12	1.97	0.47
1:A:73:TYR:O	1:A:75:LYS:HE2	2.15	0.47
1:A:112:VAL:HG23	3:A:741:HOH:O	2.13	0.47
1:A:309:PRO:C	1:A:310:ASN:HD22	2.23	0.47
1:A:455:VAL:HB	1:A:499:LEU:HD22	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:PRO:O	1:A:289:GLU:HB2	2.15	0.47
1:A:461[B]:LYS:HE3	1:A:463:ASP:OD2	2.14	0.46
1:A:647:LEU:HD12	1:A:647:LEU:O	2.15	0.46
1:A:173:PHE:HA	2:A:711:X98:C19	2.45	0.46
1:A:245:ARG:O	1:A:245:ARG:HG3	2.15	0.46
1:A:567:ARG:HH21	1:A:569:ASP:CG	2.21	0.46
1:A:245:ARG:NH2	1:A:265:ASP:OD2	2.48	0.46
1:A:310:ASN:O	1:A:311:TYR:HB2	2.16	0.46
1:A:436:GLN:HE21	1:A:438:VAL:HG11	1.80	0.46
1:A:589:TYR:CD1	1:A:647:LEU:HD21	2.50	0.46
1:A:120:SER:HB3	1:A:125:VAL:CG2	2.45	0.46
1:A:151:VAL:CG1	1:A:152:THR:N	2.79	0.46
1:A:166:ASP:OD2	1:A:213:HIS:NE2	2.36	0.46
1:A:345:ARG:HE	1:A:345:ARG:HB3	1.43	0.46
1:A:62:LEU:CD1	1:A:708:TRP:CD1	2.99	0.46
2:A:711:X98:HA1	2:A:711:X98:CB	2.47	0.45
1:A:221:GLU:O	1:A:222:ASP:C	2.59	0.45
1:A:515:HIS:O	1:A:519:ILE:HG12	2.16	0.45
1:A:151:VAL:HG12	1:A:152:THR:N	2.30	0.45
1:A:181:ASP:OD1	1:A:181:ASP:C	2.58	0.45
1:A:516:LYS:HA	1:A:519:ILE:CG1	2.46	0.45
1:A:692:GLU:O	1:A:695:ASP:HB2	2.17	0.45
1:A:450:ILE:HG21	1:A:504:ILE:HA	1.99	0.45
1:A:171:VAL:HG13	1:A:188:ASN:O	2.15	0.45
1:A:547:ARG:HA	1:A:704:LEU:CD1	2.41	0.45
1:A:380:GLY:HA2	1:A:482:PRO:HB2	1.99	0.45
1:A:173:PHE:CD2	2:A:711:X98:C4	3.00	0.45
1:A:157:LYS:HB3	1:A:162:LYS:HB2	1.99	0.44
1:A:589:TYR:O	1:A:590:THR:C	2.60	0.44
1:A:172:LYS:HG2	1:A:173:PHE:CD1	2.52	0.44
1:A:154:LYS:HE2	1:A:154:LYS:HB3	1.60	0.44
1:A:320:ASP:N	1:A:321:PRO:CD	2.81	0.44
1:A:499:LEU:HD23	1:A:499:LEU:O	2.17	0.44
1:A:71:TYR:HE2	1:A:686:THR:HG23	1.82	0.43
1:A:384:GLY:O	1:A:396:TYR:HA	2.17	0.43
1:A:534:ALA:O	1:A:535:GLU:C	2.61	0.43
1:A:493:ARG:NH1	1:A:494:HIS:HD1	2.16	0.43
1:A:310:ASN:HB2	1:A:332:GLU:OE1	2.17	0.43
1:A:576:ALA:HB3	1:A:635:LEU:HD22	1.99	0.43
1:A:57:CYS:HA	1:A:58:PRO:HD3	1.87	0.43
1:A:423:ARG:NH1	1:A:423:ARG:HG2	2.25	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:THR:HA	1:A:315:ASN:HD22	1.84	0.42
1:A:130:TYR:HB2	1:A:139:PHE:CE1	2.54	0.42
1:A:400:SER:HB3	1:A:403:SER:HB3	2.01	0.42
1:A:450:ILE:HA	1:A:451:PRO:HD3	1.83	0.42
1:A:519:ILE:CG2	1:A:598:ASP:OD2	2.66	0.42
1:A:46:GLN:O	1:A:49:ILE:HG12	2.20	0.42
1:A:100:LEU:CD2	1:A:113:PHE:HD2	2.31	0.42
1:A:154:LYS:HD2	1:A:167:VAL:HG22	2.01	0.42
1:A:623:PRO:O	1:A:665:LYS:NZ	2.52	0.42
1:A:323:GLU:HA	1:A:326:TRP:CD2	2.55	0.42
1:A:138:TYR:HE2	1:A:157:LYS:HD2	1.84	0.42
1:A:160:GLY:O	1:A:161:ALA:C	2.63	0.42
1:A:206:LEU:O	1:A:207:HIS:HB2	2.18	0.42
1:A:383:VAL:O	1:A:383:VAL:HG12	2.19	0.42
1:A:102:VAL:HG13	1:A:111:ARG:HG2	2.00	0.42
1:A:521:ALA:HB2	1:A:610:TRP:CZ2	2.55	0.42
1:A:492:VAL:HG13	1:A:497:GLY:O	2.19	0.42
1:A:647:LEU:C	1:A:647:LEU:CD1	2.83	0.41
1:A:39:THR:O	1:A:42:PHE:HB3	2.21	0.41
1:A:635:LEU:HD11	1:A:652:PHE:CD2	2.53	0.41
1:A:53:PHE:O	1:A:702:ARG:NH1	2.50	0.41
1:A:450:ILE:CG2	1:A:504:ILE:HA	2.50	0.41
1:A:166:ASP:C	1:A:167:VAL:HG23	2.45	0.41
1:A:53:PHE:CZ	1:A:669:PRO:HG3	2.55	0.41
1:A:115:ASP:OD1	1:A:117:ASN:HB2	2.21	0.41
1:A:266:LEU:HD23	1:A:266:LEU:HA	1.89	0.41
1:A:82:LYS:HG2	1:A:132:PHE:CD1	2.56	0.41
1:A:81:LYS:O	1:A:82:LYS:HD2	2.20	0.41
1:A:547:ARG:HH11	1:A:547:ARG:CG	2.33	0.41
1:A:636:LEU:CD2	1:A:673:HIS:HB3	2.51	0.41
1:A:122:ASP:OD1	1:A:122:ASP:N	2.54	0.41
1:A:395:PHE:CZ	1:A:409:HIS:CD2	3.09	0.40
1:A:140:ALA:HB2	1:A:155:PHE:CD2	2.56	0.40
1:A:554:SER:HB3	1:A:555:ASN:H	1.55	0.40
1:A:641:ASP:OD1	1:A:641:ASP:C	2.65	0.40
1:A:57:CYS:SG	1:A:59:ILE:CG2	3.04	0.40
1:A:302:PHE:CD1	1:A:316:ILE:HD12	2.56	0.40
1:A:438:VAL:O	1:A:438:VAL:CG2	2.67	0.40
1:A:563:CYS:O	1:A:571:PHE:CE1	2.70	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	710/710 (100%)	653 (92%)	51 (7%)	6 (1%)	16	29

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	346	SER
1	A	357	VAL
1	A	578	VAL
1	A	281	LYS
1	A	590	THR
1	A	474	GLY

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	605/617 (98%)	522 (86%)	83 (14%)	3	6

All (83) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	SER
1	A	13	GLU
1	A	23	LYS
1	A	24	VAL
1	A	35	ASP

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Mol	Chain	Res	Type
1	A	43	VAL
1	A	49	ILE
1	A	62	LEU
1	A	65	GLU
1	A	79	HIS
1	A	82	LYS
1	A	84	LYS
1	A	103	GLN
1	A	106	LEU
1	A	111	ARG
1	A	114	LEU
1	A	118	ILE
1	A	144	SER
1	A	163	GLU
1	A	170	ARG
1	A	183	LYS
1	A	196	LYS
1	A	204	THR
1	A	218	ASP
1	A	220	SER
1	A	230	ASP
1	A	240	LEU
1	A	241	SER
1	A	260	ARG
1	A	270	SER
1	A	271	ASN
1	A	278	LYS
1	A	281	LYS
1	A	287	GLU
1	A	294	THR
1	A	308	SER
1	A	332	GLU
1	A	335[A]	LYS
1	A	335[B]	LYS
1	A	345	ARG
1	A	347	ASN
1	A	351	LEU
1	A	357	VAL
1	A	358	LYS
1	A	359	ASN
1	A	362	GLN
1	A	368	THR

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Mol	Chain	Res	Type
1	A	381	SER
1	A	386	SER
1	A	389	LYS
1	A	393	GLU
1	A	402	LEU
1	A	414	LYS
1	A	415	GLU
1	A	423	ARG
1	A	425	VAL
1	A	428	LYS
1	A	434	ASP
1	A	438	VAL
1	A	450	ILE
1	A	479	SER
1	A	489	LEU
1	A	499	LEU
1	A	504	ILE
1	A	535	GLU
1	A	547	ARG
1	A	554	SER
1	A	567	ARG
1	A	585	LYS
1	A	598	ASP
1	A	604	SER
1	A	622	LEU
1	A	624	GLU
1	A	651	LYS
1	A	653	ILE
1	A	663	SER
1	A	664	ARG
1	A	665	LYS
1	A	676	THR
1	A	689	VAL
1	A	702	ARG
1	A	704	LEU
1	A	706	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	GLN
1	A	207	HIS

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Mol	Chain	Res	Type
1	A	310	ASN
1	A	315	ASN
1	A	347	ASN
1	A	409	HIS
1	A	436	GLN
1	A	477	ASN
1	A	551	ASN
1	A	555	ASN
1	A	577	GLN
1	A	606	GLN
1	A	673	HIS

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 ⓘ

1 ligand is 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	X98	A	711	-	37,37,37	1.09	4 (10%)	51,52,52	1.02	4 (7%)

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	X98	A	711	-	-	3/21/38/38	0/5/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	711	X98	C5-C4	2.81	1.52	1.48
2	A	711	X98	C12-C7	2.39	1.45	1.41
2	A	711	X98	C4-N3	2.25	1.42	1.38
2	A	711	X98	C11-C12	2.17	1.43	1.39

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	711	X98	C14-C5-C4	-3.17	116.05	120.44
2	A	711	X98	C14-C5-N6	2.89	121.24	117.47
2	A	711	X98	C1-C-N	-2.84	110.35	118.06
2	A	711	X98	C1-C2-N3	2.32	114.78	112.27

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	711	X98	O-C-N-CD
2	A	711	X98	C-C1-C2-N3
2	A	711	X98	O-C-N-CA

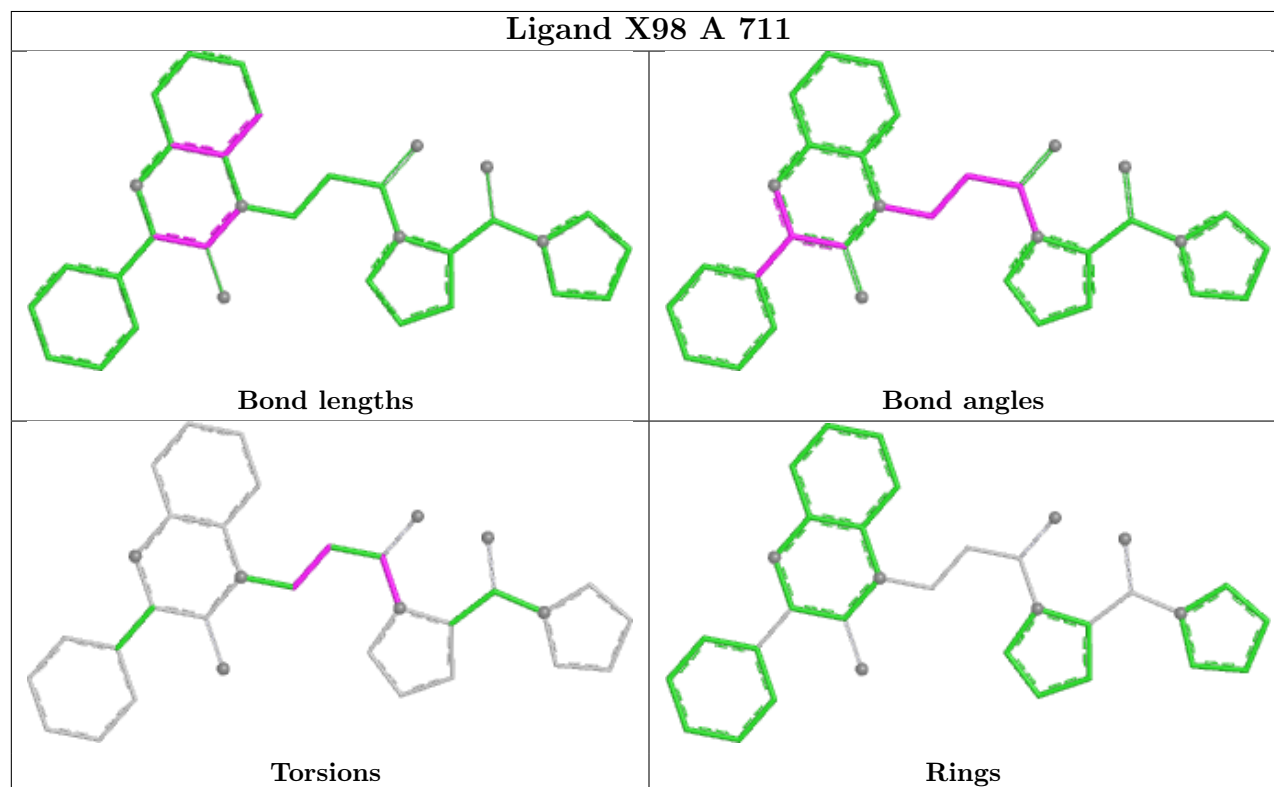
There are no ring outliers.

1 monomer is involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	711	X98	10	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 [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	709/710 (99%)	-0.15	1 (0%) 92 93	14, 32, 62, 97	3 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	LEU	2.5

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](#)

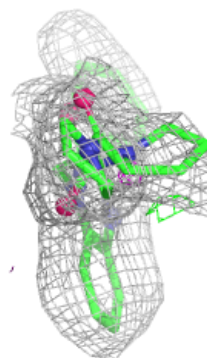
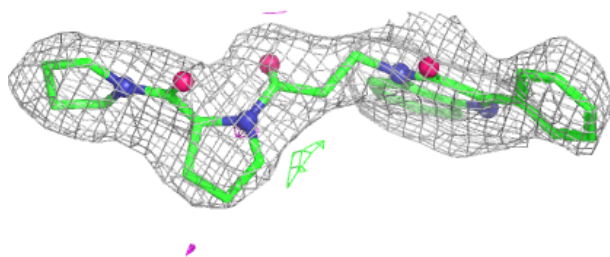
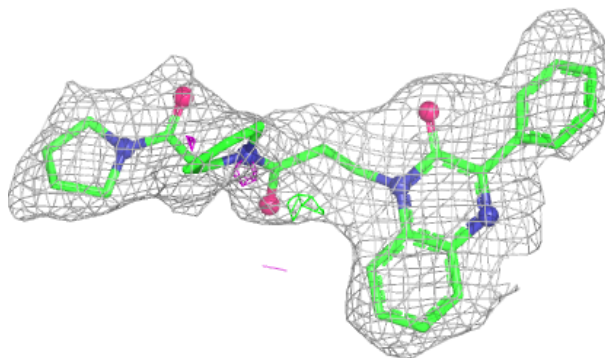
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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	X98	A	711	33/33	0.91	0.12	28,47,53,58	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 X98 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.