



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 10:38 AM UTC

PDB ID : 4TYB / pdb_00004tyb
Title : An Ligand-observed Mass Spectrometry-based Approach Integrated into the
Fragment Based Lead Discovery Pipeline
Authors : Shui, W.; Yang, C.; Lin, J.; Chen, X.; Qin, S.; Chen, S.
Deposited on : 2014-07-08
Resolution : 2.93 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
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

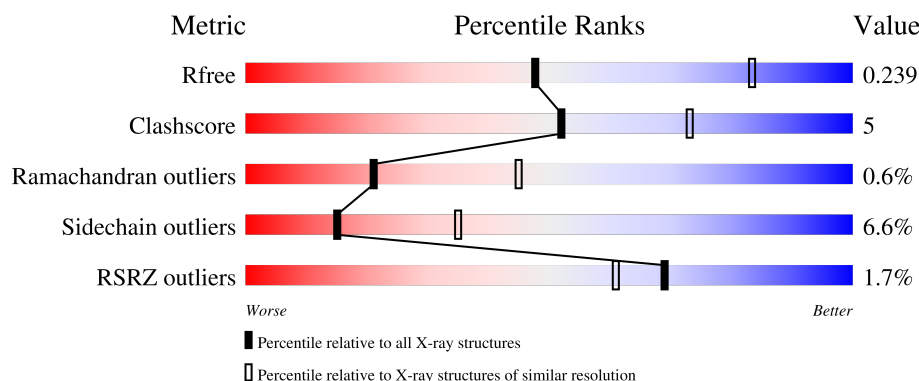
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.93 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	1159 (2.96-2.92)
Clashscore	190562	1184 (2.96-2.92)
Ramachandran outliers	187476	1131 (2.96-2.92)
Sidechain outliers	187428	1131 (2.96-2.92)
RSRZ outliers	180081	1159 (2.96-2.92)

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	566	2% 79% 15% • •
1	B	566	2% 80% 14% • •
1	C	566	• 83% 11% • •
1	D	566	• 80% 14% • •

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	3B1	A	601	-	-	X	-
2	3B1	B	601	X	-	-	-
2	3B1	C	601	X	-	X	-
2	3B1	D	601	X	-	-	-

2 Entry composition [i](#)

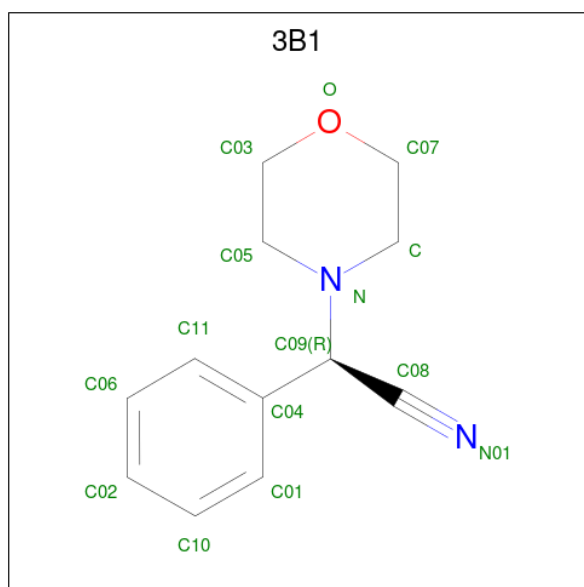
There are 2 unique types of molecules in this entry. The entry contains 16969 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	545	Total	C	N	O	S	0	0	0
			4229	2669	745	784	31			
1	B	544	Total	C	N	O	S	0	0	0
			4222	2664	744	783	31			
1	C	545	Total	C	N	O	S	0	0	0
			4229	2669	745	784	31			
1	D	545	Total	C	N	O	S	0	0	0
			4229	2669	745	784	31			

- Molecule 2 is (2R)-morpholin-4-yl(phenyl)ethanenitrile (CCD ID: 3B1) (formula: C₁₂H₁₄N₂O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			15	12	2	1		
2	B	1	Total	C	N	O	0	0
			15	12	2	1		

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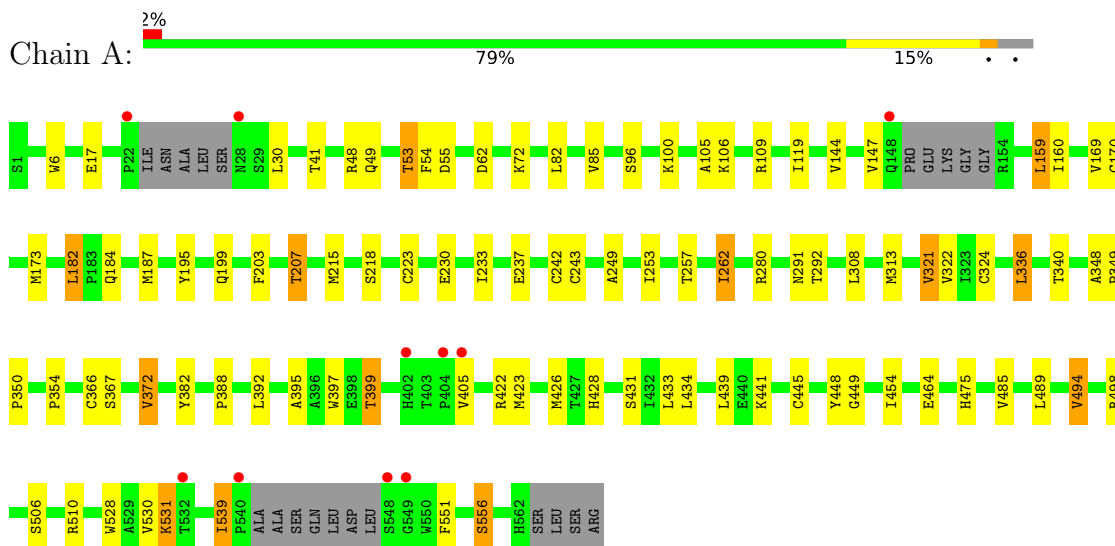
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			15	12	2	1		
2	D	1	Total	C	N	O	0	0
			15	12	2	1		

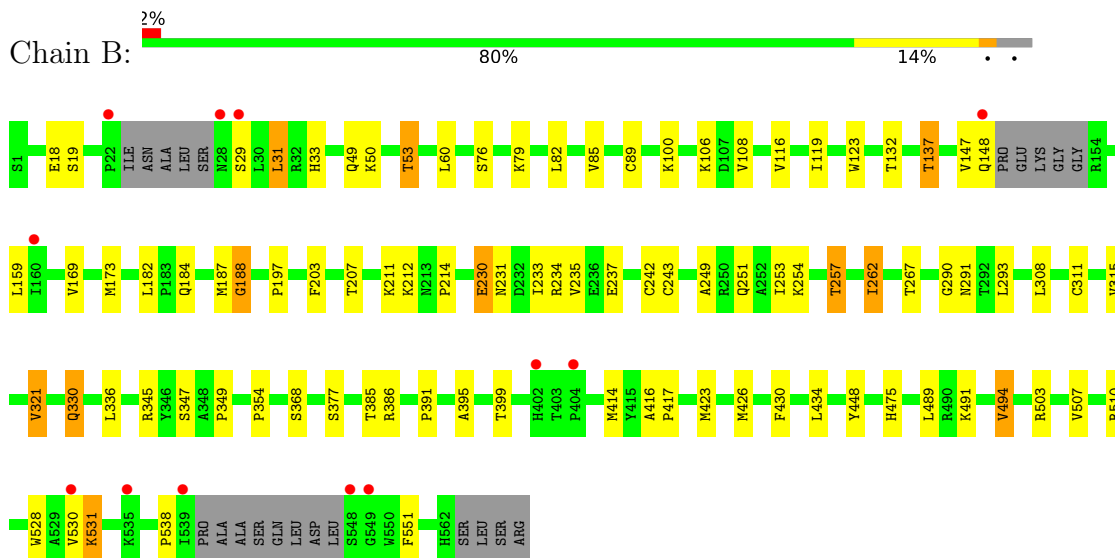
3 Residue-property plots

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.

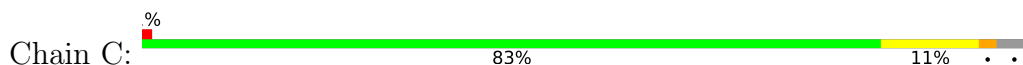
• Molecule 1: Polyprotein

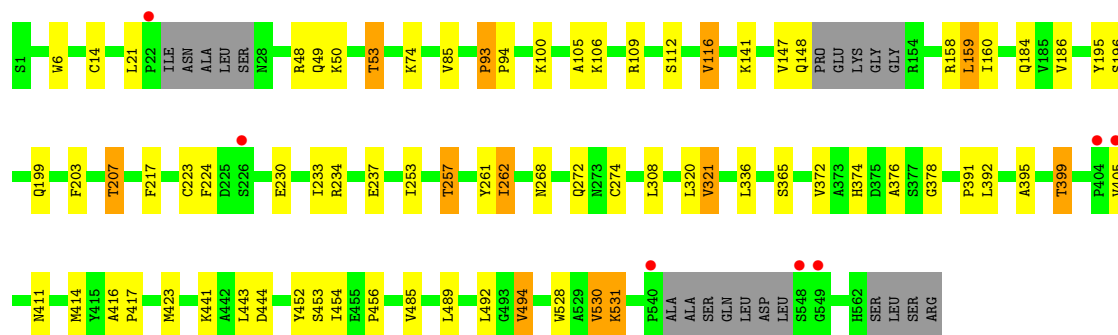


• Molecule 1: Polyprotein

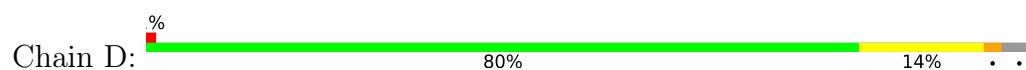


• Molecule 1: Polyprotein





• Molecule 1: Polyprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	102.45Å 102.35Å 251.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.41 – 2.93 47.41 – 2.93	Depositor EDS
% Data completeness (in resolution range)	98.6 (47.41-2.93) 98.8 (47.41-2.93)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.69 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.8.0049, REFMAC 5.8.0049	Depositor
R, R_{free}	0.205 , 0.273 (Not available) , 0.239	Depositor DCC
R_{free} test set	2899 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	33.9	Xtriage
Anisotropy	0.065	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 10.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.087 for k,h,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	16969	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.48% 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: 3B1

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	0.74	0/4321	0.95	3/5864 (0.1%)
1	B	0.73	0/4313	0.98	5/5852 (0.1%)
1	C	0.73	0/4321	0.97	4/5864 (0.1%)
1	D	0.74	0/4321	0.96	8/5864 (0.1%)
All	All	0.74	0/17276	0.97	20/23444 (0.1%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	321	VAL	CB-CA-C	-6.82	100.00	110.50
1	A	321	VAL	CB-CA-C	-6.17	101.00	110.50
1	B	349	PRO	O-C-N	5.97	123.96	121.15
1	D	530	VAL	N-CA-C	5.88	116.92	111.45
1	B	345	ARG	N-CA-C	-5.80	104.86	111.07
1	D	321	VAL	CB-CA-C	-5.57	101.07	110.71
1	C	93	PRO	O-C-N	5.53	123.75	121.15
1	D	352	ASP	CA-C-N	-5.53	114.69	120.38
1	D	352	ASP	C-N-CA	-5.53	114.69	120.38
1	D	475	HIS	N-CA-C	5.47	122.45	110.80
1	B	475	HIS	CA-C-N	-5.36	114.99	123.24
1	B	475	HIS	C-N-CA	-5.36	114.99	123.24
1	A	218	SER	N-CA-C	-5.12	101.92	110.17
1	C	186	VAL	N-CA-C	-5.08	105.64	110.42
1	B	321	VAL	CB-CA-C	-5.06	102.29	110.28
1	D	192	GLY	N-CA-C	5.05	120.27	114.16
1	D	475	HIS	CA-C-N	-5.02	116.27	123.20
1	D	475	HIS	C-N-CA	-5.02	116.27	123.20
1	C	224	PHE	N-CA-C	5.02	116.75	111.28
1	A	475	HIS	N-CA-C	5.02	121.59	114.16

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	4229	0	4219	48	0
1	B	4222	0	4212	40	0
1	C	4229	0	4219	42	0
1	D	4229	0	4219	38	0
2	A	15	0	14	10	0
2	B	15	0	14	1	0
2	C	15	0	14	6	0
2	D	15	0	14	0	0
All	All	16969	0	16925	174	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 5.

All (174) 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:B:49:GLN:O	1:B:53:THR:HG22	1.55	1.06
1:D:49:GLN:O	1:D:53:THR:HG22	1.54	1.05
1:A:49:GLN:O	1:A:53:THR:HG23	1.62	0.99
1:A:366:CYS:HB3	2:A:601:3B1:H14	1.45	0.98
1:A:395:ALA:O	1:A:399:THR:HG22	1.72	0.89
1:D:203:PHE:O	1:D:207:THR:HG23	1.73	0.88
1:B:49:GLN:O	1:B:53:THR:CG2	2.26	0.83
1:C:253:ILE:O	1:C:257:THR:HG23	1.79	0.83
1:A:6:TRP:HE1	1:A:53:THR:HG21	1.43	0.81
1:A:366:CYS:HB3	2:A:601:3B1:C	2.11	0.79
1:C:203:PHE:O	1:C:207:THR:HG23	1.83	0.78
1:A:203:PHE:O	1:A:207:THR:HG23	1.83	0.78
1:B:395:ALA:O	1:B:399:THR:HG23	1.83	0.77
1:A:253:ILE:O	1:A:257:THR:HG23	1.85	0.77
1:C:49:GLN:O	1:C:53:THR:CG2	2.33	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:82:LEU:HD13	1:D:249:ALA:HB2	1.67	0.76
1:C:49:GLN:O	1:C:53:THR:HG23	1.89	0.72
1:D:49:GLN:O	1:D:53:THR:CG2	2.38	0.69
1:C:234:ARG:HG3	1:C:262:ILE:HD11	1.76	0.66
1:B:123:TRP:CE3	1:B:173:MET:HE3	2.32	0.65
1:D:253:ILE:O	1:D:257:THR:HG23	1.97	0.64
1:A:366:CYS:SG	2:A:601:3B1:H11	2.38	0.64
1:A:340:THR:HG23	1:A:350:PRO:HG3	1.81	0.63
1:A:48:ARG:HG2	1:A:159:LEU:HD13	1.80	0.62
1:B:231:ASN:O	1:B:235:VAL:HG23	2.00	0.62
1:D:395:ALA:O	1:D:399:THR:HG22	2.02	0.60
2:A:601:3B1:C	2:A:601:3B1:H2	2.31	0.60
1:A:82:LEU:HD13	1:A:249:ALA:HB2	1.86	0.58
1:B:182:LEU:HD13	1:B:243:CYS:SG	2.43	0.58
2:A:601:3B1:H2	2:A:601:3B1:H15	1.85	0.57
1:A:366:CYS:CB	2:A:601:3B1:H14	2.26	0.57
1:B:233:ILE:HG22	1:B:262:ILE:HD12	1.87	0.57
1:D:31:LEU:O	1:D:31:LEU:HD12	2.06	0.56
1:A:449:GLY:HA3	1:A:556:SER:CA	2.36	0.55
1:A:423:MET:HE3	1:A:528:TRP:CZ3	2.41	0.55
1:D:308:LEU:HB3	1:D:311:CYS:SG	2.46	0.55
1:A:105:ALA:O	1:A:109:ARG:HG3	2.06	0.55
1:A:144:VAL:HG21	1:A:397:TRP:CG	2.42	0.55
1:B:182:LEU:CD1	1:B:243:CYS:SG	2.95	0.55
1:A:336:LEU:HD12	1:A:354:PRO:HG2	1.87	0.54
1:C:49:GLN:O	1:C:53:THR:HG22	2.07	0.54
1:A:237:GLU:HG3	1:A:257:THR:HG21	1.90	0.54
1:A:449:GLY:HA3	1:A:556:SER:HA	1.90	0.54
1:B:423:MET:HA	1:B:528:TRP:CZ2	2.43	0.54
1:D:263:GLY:HA3	1:D:277:ARG:O	2.08	0.54
1:B:253:ILE:O	1:B:257:THR:HG23	2.08	0.54
1:D:241:GLN:OE1	1:D:250:ARG:HG3	2.08	0.53
1:D:530:VAL:O	1:D:531:LYS:CB	2.57	0.53
2:A:601:3B1:H2	2:A:601:3B1:C07	2.39	0.53
1:B:503:ARG:O	1:B:507:VAL:HG23	2.09	0.53
1:A:489:LEU:HA	1:A:494:VAL:HG13	1.91	0.53
1:D:30:LEU:O	1:D:494:VAL:HB	2.09	0.52
1:D:237:GLU:HG3	1:D:257:THR:HG21	1.92	0.52
1:B:233:ILE:CG2	1:B:262:ILE:HD12	2.40	0.52
1:C:321:VAL:CG2	1:C:365:SER:CB	2.88	0.51
1:A:6:TRP:NE1	1:A:53:THR:HG21	2.21	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:VAL:HG11	1:B:116:VAL:HG13	1.91	0.51
1:B:434:LEU:HD21	1:B:510:ARG:HB2	1.91	0.51
1:B:203:PHE:O	1:B:207:THR:HG23	2.10	0.51
1:B:530:VAL:O	1:B:531:LYS:CB	2.59	0.50
1:C:147:VAL:HG12	1:C:148:GLN:HG2	1.93	0.50
1:A:464:GLU:OE2	1:A:539:ILE:HG12	2.10	0.50
1:B:187:MET:O	1:B:188:GLY:C	2.54	0.50
1:B:426:MET:O	1:B:430:PHE:HB2	2.11	0.49
1:C:530:VAL:O	1:C:531:LYS:HB3	2.12	0.49
1:D:385:THR:OG1	1:D:386:ARG:N	2.46	0.49
1:A:280:ARG:NE	1:A:291:ASN:OD1	2.36	0.49
1:B:237:GLU:CG	1:B:257:THR:HG21	2.42	0.49
1:B:33:HIS:HD2	1:B:491:LYS:O	1.96	0.49
1:C:416:ALA:HB3	1:C:417:PRO:HD3	1.95	0.49
1:D:119:ILE:HD13	1:D:169:VAL:HG11	1.94	0.49
1:B:197:PRO:HG3	1:B:414:MET:CE	2.43	0.48
1:B:489:LEU:HD22	1:B:494:VAL:HG22	1.95	0.48
1:A:313:MET:HG2	1:A:322:VAL:HG22	1.95	0.48
1:D:170:CYS:HA	1:D:173:MET:HE3	1.95	0.48
1:C:414:MET:HE2	2:C:601:3B1:H5	1.95	0.48
1:D:17:GLU:OE1	1:D:41:THR:HB	2.13	0.48
1:A:434:LEU:HD21	1:A:510:ARG:CB	2.43	0.48
2:A:601:3B1:H2	2:A:601:3B1:H10	1.96	0.48
1:C:391:PRO:HB3	1:C:411:ASN:ND2	2.29	0.48
1:C:414:MET:HG3	2:C:601:3B1:C03	2.44	0.47
2:A:601:3B1:H15	2:A:601:3B1:C01	2.44	0.47
1:B:76:SER:HA	1:B:242:CYS:O	2.14	0.47
1:D:48:ARG:HG2	1:D:159:LEU:HD13	1.94	0.47
1:C:85:VAL:HG11	1:C:116:VAL:HG22	1.96	0.47
1:C:195:TYR:HA	1:C:199:GLN:HE21	1.79	0.47
1:C:6:TRP:HE1	1:C:53:THR:HG21	1.79	0.47
1:A:54:PHE:HD1	1:A:55:ASP:O	1.98	0.46
1:C:321:VAL:HG22	1:C:365:SER:HB3	1.98	0.46
1:C:423:MET:HA	1:C:528:TRP:CZ2	2.50	0.46
1:A:233:ILE:CG2	1:A:262:ILE:HD12	2.46	0.46
1:A:422:ARG:HA	1:A:426:MET:SD	2.55	0.46
1:D:515:GLY:HA2	1:D:519:ALA:HB2	1.97	0.46
1:B:187:MET:HE2	1:B:293:LEU:HD23	1.96	0.46
1:B:330:GLN:HE21	1:B:330:GLN:HA	1.80	0.46
1:B:414:MET:HE2	2:B:601:3B1:H12	1.98	0.46
1:D:196:SER:H	1:D:199:GLN:HE21	1.63	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:40:THR:HB	1:D:157:ALA:HB2	1.96	0.46
1:B:336:LEU:HD12	1:B:354:PRO:HB2	1.98	0.46
1:A:372:VAL:HG12	1:A:382:TYR:CD1	2.50	0.45
1:A:530:VAL:O	1:A:531:LYS:CB	2.64	0.45
1:C:233:ILE:CG2	1:C:262:ILE:HD12	2.46	0.45
1:D:196:SER:O	1:D:197:PRO:C	2.56	0.45
1:A:434:LEU:HD21	1:A:510:ARG:HB2	1.98	0.45
1:C:374:HIS:HD2	1:C:378:GLY:O	1.99	0.45
1:A:448:TYR:CE2	1:A:551:PHE:HD1	2.34	0.45
1:C:485:VAL:O	1:C:489:LEU:HG	2.16	0.45
1:C:444:ASP:HA	1:C:452:TYR:O	2.17	0.45
1:A:428:HIS:O	1:A:431:SER:HB3	2.17	0.44
1:D:5:THR:O	1:D:275:GLY:HA3	2.17	0.44
1:D:141:LYS:HE3	1:D:143:GLU:HG3	1.99	0.44
1:A:170:CYS:HA	1:A:173:MET:HE3	1.99	0.44
1:A:348:ALA:N	1:A:349:PRO:CD	2.80	0.44
1:C:48:ARG:HG2	1:C:159:LEU:HD13	2.00	0.44
1:D:489:LEU:HA	1:D:494:VAL:HG13	1.99	0.44
1:D:268:ASN:OD1	1:D:268:ASN:C	2.61	0.44
1:A:17:GLU:HB3	1:A:41:THR:HG22	2.00	0.44
1:B:251:GLN:OE1	1:B:251:GLN:HA	2.17	0.44
1:B:385:THR:OG1	1:B:386:ARG:N	2.50	0.44
1:D:6:TRP:HE1	1:D:53:THR:HG21	1.83	0.44
1:B:448:TYR:CE2	1:B:551:PHE:HD1	2.36	0.43
1:C:395:ALA:O	1:C:399:THR:HG23	2.17	0.43
1:C:257:THR:HA	1:C:261:TYR:HB2	1.99	0.43
1:D:448:TYR:CE2	1:D:551:PHE:HD1	2.36	0.43
1:C:443:LEU:HD12	1:C:456:PRO:HG3	1.99	0.43
1:B:230:GLU:O	1:B:234:ARG:HG3	2.19	0.43
1:A:187:MET:HE1	1:A:292:THR:HG22	2.00	0.43
1:A:215:MET:HE3	1:A:324:CYS:SG	2.59	0.43
1:A:388:PRO:O	1:A:392:LEU:HG	2.17	0.43
1:B:137:THR:HA	1:B:267:THR:O	2.19	0.43
1:C:237:GLU:CG	1:C:257:THR:HG21	2.49	0.43
1:D:423:MET:HA	1:D:528:TRP:CZ2	2.54	0.43
1:B:31:LEU:HD12	1:B:31:LEU:C	2.44	0.43
1:C:217:PHE:CD1	1:C:336:LEU:HD21	2.53	0.43
1:B:290:GLY:O	1:B:291:ASN:C	2.62	0.42
1:A:433:LEU:HB3	1:A:439:LEU:HD23	2.00	0.42
1:A:96:SER:HB3	1:A:105:ALA:HB2	2.01	0.42
1:B:416:ALA:HB3	1:B:417:PRO:HD3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:414:MET:HG3	2:C:601:3B1:H5	2.01	0.42
1:D:219:TYR:HB3	1:D:320:LEU:HD23	2.01	0.42
1:D:33:HIS:HD2	1:D:491:LYS:O	2.02	0.42
1:A:119:ILE:HD13	1:A:169:VAL:HG11	2.02	0.42
1:C:489:LEU:HD22	1:C:494:VAL:CG2	2.49	0.42
1:A:72:LYS:HG2	1:A:242:CYS:SG	2.60	0.42
1:C:321:VAL:CG2	1:C:365:SER:HB2	2.49	0.42
1:D:182:LEU:N	1:D:183:PRO:CD	2.82	0.42
1:D:175:LEU:HG	1:D:253:ILE:HG12	2.01	0.42
1:C:48:ARG:HG3	1:C:159:LEU:HD22	2.02	0.42
1:D:263:GLY:CA	1:D:277:ARG:O	2.67	0.42
2:A:601:3B1:C	2:A:601:3B1:C01	2.96	0.42
1:B:530:VAL:O	1:B:531:LYS:HB3	2.20	0.42
1:D:89:CYS:SG	1:D:108:VAL:HG13	2.59	0.42
1:C:268:ASN:HB3	1:C:274:CYS:SG	2.60	0.42
1:D:195:TYR:HA	1:D:199:GLN:HE21	1.85	0.42
1:A:30:LEU:O	1:A:494:VAL:HB	2.20	0.41
1:A:182:LEU:HD13	1:A:243:CYS:SG	2.60	0.41
1:C:321:VAL:HG22	1:C:365:SER:CB	2.49	0.41
1:C:414:MET:CE	2:C:601:3B1:H6	2.51	0.41
1:D:31:LEU:HD12	1:D:31:LEU:C	2.45	0.41
1:B:119:ILE:HD13	1:B:169:VAL:HG11	2.02	0.41
1:D:423:MET:HE3	1:D:528:TRP:CZ3	2.56	0.41
1:A:485:VAL:O	1:A:489:LEU:HG	2.21	0.41
1:C:105:ALA:O	1:C:109:ARG:HG3	2.20	0.41
1:A:195:TYR:HA	1:A:199:GLN:HE21	1.86	0.41
1:C:237:GLU:HG3	1:C:257:THR:HG21	2.01	0.41
1:C:321:VAL:CG2	1:C:365:SER:HB3	2.51	0.41
1:C:392:LEU:HD23	1:C:392:LEU:HA	1.94	0.41
1:C:414:MET:HE2	2:C:601:3B1:H6	2.02	0.41
1:B:211:LYS:HB2	1:B:214:PRO:HB3	2.02	0.40
1:A:445:CYS:SG	1:A:454:ILE:HD12	2.61	0.40
1:C:414:MET:HE2	2:C:601:3B1:H12	2.03	0.40
1:B:89:CYS:SG	1:B:108:VAL:HG13	2.62	0.40
1:C:93:PRO:HA	1:C:94:PRO:HD3	1.94	0.40
1:B:82:LEU:HD13	1:B:249:ALA:HB2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	537/566 (95%)	505 (94%)	29 (5%)	3 (1%)	21	45
1	B	536/566 (95%)	506 (94%)	26 (5%)	4 (1%)	18	40
1	C	537/566 (95%)	510 (95%)	24 (4%)	3 (1%)	21	45
1	D	537/566 (95%)	506 (94%)	29 (5%)	2 (0%)	30	54
All	All	2147/2264 (95%)	2027 (94%)	108 (5%)	12 (1%)	21	45

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	531	LYS
1	A	539	ILE
1	B	531	LYS
1	D	531	LYS
1	D	230	GLU
1	A	223	CYS
1	B	538	PRO
1	B	188	GLY
1	B	347	SER
1	C	223	CYS
1	C	376	ALA
1	C	531	LYS

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	461/482 (96%)	436 (95%)	25 (5%)	20	42
1	B	460/482 (95%)	430 (94%)	30 (6%)	15	35
1	C	461/482 (96%)	430 (93%)	31 (7%)	15	34
1	D	461/482 (96%)	426 (92%)	35 (8%)	12	29
All	All	1843/1928 (96%)	1722 (93%)	121 (7%)	15	35

All (121) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	THR
1	A	62	ASP
1	A	85	VAL
1	A	100	LYS
1	A	106	LYS
1	A	147	VAL
1	A	159	LEU
1	A	160	ILE
1	A	182	LEU
1	A	184	GLN
1	A	207	THR
1	A	230	GLU
1	A	262	ILE
1	A	308	LEU
1	A	321	VAL
1	A	336	LEU
1	A	367	SER
1	A	372	VAL
1	A	399	THR
1	A	405	VAL
1	A	441	LYS
1	A	494	VAL
1	A	498	ARG
1	A	506	SER
1	A	556	SER
1	B	18	GLU
1	B	19	SER
1	B	29	SER
1	B	31	LEU
1	B	50	LYS
1	B	53	THR
1	B	60	LEU
1	B	79	LYS

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Mol	Chain	Res	Type
1	B	100	LYS
1	B	106	LYS
1	B	132	THR
1	B	137	THR
1	B	147	VAL
1	B	148	GLN
1	B	159	LEU
1	B	184	GLN
1	B	212	LYS
1	B	230	GLU
1	B	254	LYS
1	B	257	THR
1	B	262	ILE
1	B	308	LEU
1	B	311	CYS
1	B	315	VAL
1	B	321	VAL
1	B	330	GLN
1	B	368	SER
1	B	377	SER
1	B	391	PRO
1	B	494	VAL
1	C	14	CYS
1	C	21	LEU
1	C	50	LYS
1	C	53	THR
1	C	74	LYS
1	C	100	LYS
1	C	106	LYS
1	C	112	SER
1	C	116	VAL
1	C	141	LYS
1	C	158	ARG
1	C	159	LEU
1	C	160	ILE
1	C	184	GLN
1	C	196	SER
1	C	207	THR
1	C	230	GLU
1	C	257	THR
1	C	262	ILE
1	C	272	GLN

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Mol	Chain	Res	Type
1	C	308	LEU
1	C	320	LEU
1	C	372	VAL
1	C	399	THR
1	C	405	VAL
1	C	441	LYS
1	C	453	SER
1	C	454	ILE
1	C	492	LEU
1	C	494	VAL
1	C	530	VAL
1	D	18	GLU
1	D	31	LEU
1	D	51	LYS
1	D	53	THR
1	D	100	LYS
1	D	106	LYS
1	D	112	SER
1	D	130	THR
1	D	158	ARG
1	D	159	LEU
1	D	160	ILE
1	D	207	THR
1	D	226	SER
1	D	230	GLU
1	D	257	THR
1	D	262	ILE
1	D	321	VAL
1	D	335	SER
1	D	336	LEU
1	D	355	GLN
1	D	361	GLU
1	D	371	SER
1	D	377	SER
1	D	399	THR
1	D	405	VAL
1	D	432	ILE
1	D	434	LEU
1	D	441	LYS
1	D	476	SER
1	D	487	SER
1	D	494	VAL

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Mol	Chain	Res	Type
1	D	498	ARG
1	D	508	ARG
1	D	534	LEU
1	D	539	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	HIS
1	A	58	GLN
1	A	199	GLN
1	A	273	ASN
1	A	467	HIS
1	A	562	HIS
1	B	28	ASN
1	B	33	HIS
1	B	58	GLN
1	B	120	HIS
1	B	199	GLN
1	B	330	GLN
1	B	514	GLN
1	B	562	HIS
1	C	33	HIS
1	C	120	HIS
1	C	199	GLN
1	C	231	ASN
1	C	374	HIS
1	C	402	HIS
1	D	33	HIS
1	D	58	GLN
1	D	120	HIS
1	D	199	GLN
1	D	355	GLN
1	D	374	HIS
1	D	562	HIS

5.3.3 RNA ⓘ

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](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	3B1	A	601	-	15,16,16	1.52	1 (6%)	15,20,20	1.64	3 (20%)
2	3B1	D	601	-	15,16,16	1.54	2 (13%)	15,20,20	1.33	2 (13%)
2	3B1	C	601	-	15,16,16	1.86	4 (26%)	15,20,20	2.17	5 (33%)
2	3B1	B	601	-	15,16,16	1.84	3 (20%)	15,20,20	3.20	5 (33%)

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	3B1	D	601	-	1/1/2/2	4/8/18/18	0/2/2/2
2	3B1	A	601	-	-	3/8/18/18	0/2/2/2
2	3B1	C	601	-	1/1/2/2	4/8/18/18	0/2/2/2
2	3B1	B	601	-	1/1/2/2	3/8/18/18	0/2/2/2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	3B1	C04-C09	4.80	1.57	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	3B1	C04-C09	4.69	1.57	1.52
2	A	601	3B1	C08-N01	3.78	1.21	1.14
2	D	601	3B1	C04-C09	3.26	1.55	1.52
2	C	601	3B1	C11-C04	3.19	1.44	1.39
2	B	601	3B1	C11-C04	3.06	1.44	1.39
2	D	601	3B1	C11-C04	2.74	1.43	1.39
2	C	601	3B1	C05-N	-2.45	1.42	1.47
2	C	601	3B1	C02-C10	2.34	1.43	1.38
2	B	601	3B1	C08-N01	2.22	1.18	1.14

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	3B1	C09-C08-N01	-9.94	162.13	177.51
2	B	601	3B1	C03-C05-N	5.11	114.72	109.69
2	C	601	3B1	C09-C08-N01	-4.87	169.98	177.51
2	A	601	3B1	C03-C05-N	4.01	113.64	109.69
2	C	601	3B1	C07-O-C03	3.58	121.46	109.88
2	C	601	3B1	C07-C-N	3.47	113.11	109.69
2	B	601	3B1	C07-O-C03	3.40	120.89	109.88
2	D	601	3B1	C09-C08-N01	-3.09	172.73	177.51
2	C	601	3B1	O-C03-C05	2.77	117.74	111.77
2	A	601	3B1	C-N-C05	-2.65	104.08	109.13
2	B	601	3B1	C-N-C05	-2.49	104.38	109.13
2	D	601	3B1	C03-C05-N	2.40	112.06	109.69
2	B	601	3B1	O-C03-C05	2.39	116.91	111.77
2	C	601	3B1	C-N-C05	-2.37	104.62	109.13
2	A	601	3B1	C09-C08-N01	-2.36	173.85	177.51

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	601	3B1	C09
2	C	601	3B1	C09
2	D	601	3B1	C09

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	3B1	C01-C04-C09-C08
2	A	601	3B1	C04-C09-N-C
2	A	601	3B1	C08-C09-N-C

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Mol	Chain	Res	Type	Atoms
2	B	601	3B1	C04-C09-N-C
2	B	601	3B1	C08-C09-N-C
2	C	601	3B1	C04-C09-N-C05
2	C	601	3B1	C08-C09-N-C05
2	C	601	3B1	C04-C09-N-C
2	D	601	3B1	C01-C04-C09-C08
2	D	601	3B1	C04-C09-N-C05
2	D	601	3B1	C08-C09-N-C05
2	D	601	3B1	C04-C09-N-C
2	B	601	3B1	C04-C09-N-C05
2	C	601	3B1	C08-C09-N-C

There are no ring outliers.

3 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	3B1	10	0
2	C	601	3B1	6	0
2	B	601	3B1	1	0

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	545/566 (96%)	-0.05	10 (1%) 67 59	16, 27, 45, 69	0
1	B	544/566 (96%)	-0.03	12 (2%) 62 53	14, 27, 47, 82	0
1	C	545/566 (96%)	-0.07	7 (1%) 75 68	15, 26, 45, 81	0
1	D	545/566 (96%)	-0.05	8 (1%) 72 65	14, 26, 47, 75	0
All	All	2179/2264 (96%)	-0.05	37 (1%) 69 61	14, 26, 46, 82	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	540	PRO	4.1
1	B	148	GLN	3.5
1	B	539	ILE	3.4
1	A	148	GLN	3.3
1	B	530	VAL	3.3
1	C	405	VAL	3.3
1	D	557	GLY	3.2
1	B	404	PRO	3.2
1	D	548	SER	3.1
1	B	22	PRO	3.0
1	B	548	SER	3.0
1	D	148	GLN	2.9
1	A	548	SER	2.9
1	C	404	PRO	2.8
1	A	540	PRO	2.8
1	C	22	PRO	2.8
1	A	405	VAL	2.7
1	B	549	GLY	2.6
1	C	540	PRO	2.6
1	A	404	PRO	2.5
1	A	402	HIS	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	402	HIS	2.5
1	C	549	GLY	2.5
1	C	226	SER	2.4
1	B	28	ASN	2.3
1	A	549	GLY	2.3
1	D	402	HIS	2.3
1	D	22	PRO	2.3
1	A	28	ASN	2.2
1	A	22	PRO	2.2
1	D	226	SER	2.2
1	D	405	VAL	2.1
1	B	29	SER	2.1
1	B	160	ILE	2.1
1	C	548	SER	2.1
1	B	535	LYS	2.1
1	A	532	THR	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	3B1	B	601	15/15	0.82	0.21	46,49,53,54	0
2	3B1	C	601	15/15	0.86	0.17	33,36,39,39	0
2	3B1	D	601	15/15	0.86	0.20	44,48,52,53	0
2	3B1	A	601	15/15	0.87	0.21	42,56,66,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.