



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:04 AM GMT

PDB ID : 1ZGS
Title : Parkia platycephala seed lectin in complex with 5-bromo-4-chloro-3-indolyl-a-D-mannose
Authors : Gallego del Sol, F.; Cavada, B.S.; Calvete, J.J.
Deposited on : 2005-04-22
Resolution : 2.50 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

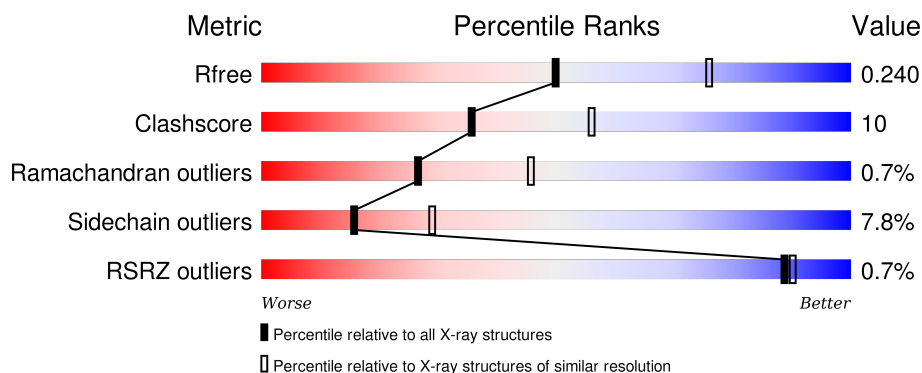
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	447	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 5%, orange 15%, yellow 21%, green 75%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 75% 21% ... </div> </div>
1	B	447	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 5%, orange 15%, yellow 17%, green 78%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 78% 17% ... </div> </div>

2 Entry composition [i](#)

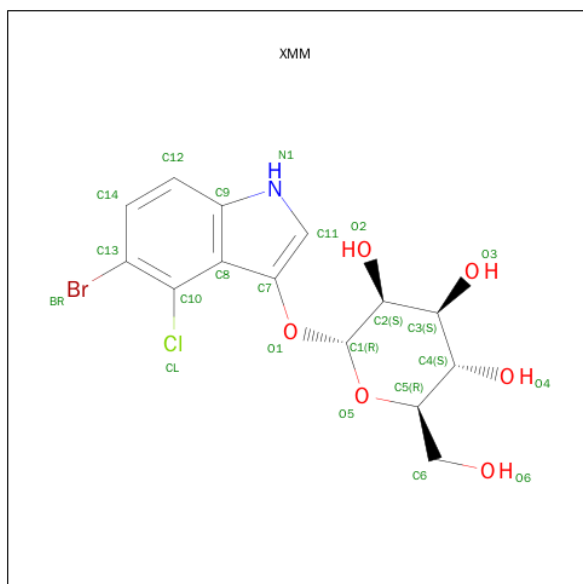
There are 3 unique types of molecules in this entry. The entry contains 7009 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Mannose/glucose-specific lectin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	442	Total	C	N	O	S	36	0	0
			3334	2142	543	646	3			
1	B	442	Total	C	N	O	S	30	0	0
			3334	2142	543	646	3			

- Molecule 2 is SUGAR ((2R,3S,4S,5S,6R)-2-(5-BROMO-4-CHLORO-1H-INDOL-3-YLOXY)-TETRAHYDRO-6-(HYDROXYMETHYL)-2H-PYRAN-3,4,5-TRIOX) (three-letter code: XMM) (formula: C₁₄H₁₅BrClNO₆).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 23	Br 1	C 14	Cl 1	N 1	O 6	0	0
2	A	1	Total 23	Br 1	C 14	Cl 1	N 1	O 6	0	0
2	A	1	Total 23	Br 1	C 14	Cl 1	N 1	O 6	0	0

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	B	1	Total 23	Br 1	C 14	Cl 1	N 1	O 6	0	0
2	B	1	Total 23	Br 1	C 14	Cl 1	N 1	O 6	0	0
2	B	1	Total 23	Br 1	C 14	Cl 1	N 1	O 6	0	0

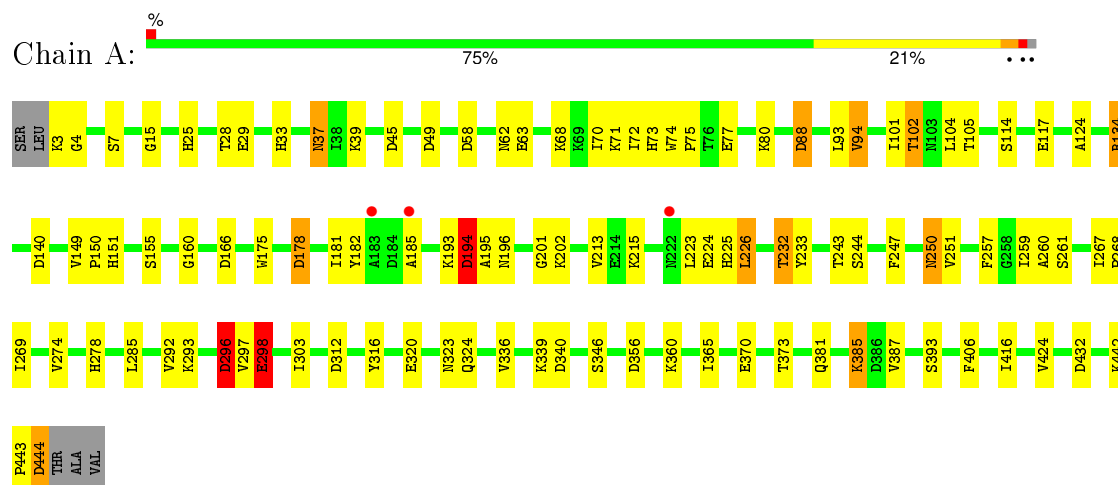
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	103	Total 103	O 103	0	0
3	B	100	Total 100	O 100	0	0

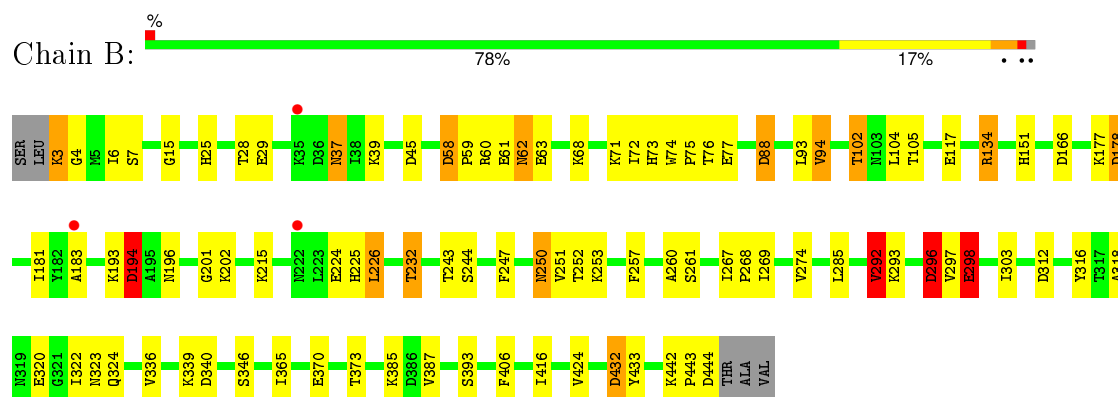
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Mannose/glucose-specific lectin



- Molecule 1: Mannose/glucose-specific lectin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	80.24Å 114.18Å 80.30Å 90.00° 119.88° 90.00°	Depositor
Resolution (Å)	30.00 – 2.50 69.63 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (30.00-2.50) 98.1 (69.63-2.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.38 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
R, R_{free}	0.208 , 0.244 0.211 , 0.240	Depositor DCC
R_{free} test set	3930 reflections (10.13%)	DCC
Wilson B-factor (Å ²)	29.6	Xtriage
Anisotropy	0.167	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 2.1	EDS
Estimated twinning fraction	0.014 for -h-l,k,h 0.014 for l,k,-h-l 0.015 for h,-k,-h-l 0.015 for -h-l,-k,l 0.487 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 43418 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7009	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.15% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: XMM

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	1.27	6/3428 (0.2%)	0.95	18/4650 (0.4%)
1	B	1.27	5/3428 (0.1%)	0.97	18/4650 (0.4%)
All	All	1.27	11/6856 (0.2%)	0.96	36/9300 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	3
All	All	0	6

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	298	GLU	C-N	-56.91	0.30	1.33
1	B	298	GLU	C-N	-55.75	0.32	1.33
1	B	296	ASP	C-N	-18.21	0.92	1.34
1	A	296	ASP	C-N	-18.19	0.92	1.34
1	B	292	VAL	CB-CG2	-7.95	1.36	1.52
1	A	3	LYS	C-N	-7.55	1.19	1.33
1	B	292	VAL	CB-CG1	-6.94	1.38	1.52
1	B	3	LYS	C-N	-6.36	1.21	1.33
1	A	292	VAL	CB-CG1	-6.27	1.39	1.52
1	A	381	GLN	CG-CD	5.92	1.64	1.51
1	A	292	VAL	CB-CG2	-5.56	1.41	1.52

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	3	LYS	O-C-N	-11.56	103.55	123.20
1	B	3	LYS	O-C-N	-11.53	103.60	123.20
1	B	3	LYS	CA-C-N	11.20	138.59	116.20
1	A	3	LYS	CA-C-N	11.08	138.37	116.20
1	A	3	LYS	C-N-CA	10.37	144.07	122.30
1	B	3	LYS	C-N-CA	10.01	143.31	122.30
1	B	202	LYS	N-CA-C	-9.85	84.41	111.00
1	A	202	LYS	N-CA-C	-9.53	85.27	111.00
1	A	194	ASP	CB-CG-OD2	8.21	125.69	118.30
1	B	194	ASP	CB-CG-OD2	7.71	125.23	118.30
1	A	223	LEU	CB-CG-CD2	7.00	122.90	111.00
1	B	183	ALA	O-C-N	6.60	133.26	122.70
1	A	297	VAL	N-CA-C	-6.07	94.62	111.00
1	A	381	GLN	CB-CG-CD	-6.03	95.92	111.60
1	B	88	ASP	CB-CG-OD2	6.00	123.70	118.30
1	A	88	ASP	CB-CG-OD2	5.93	123.64	118.30
1	A	292	VAL	CG1-CB-CG2	-5.86	101.53	110.90
1	A	432	ASP	CB-CG-OD2	5.77	123.49	118.30
1	B	296	ASP	O-C-N	5.75	131.90	122.70
1	A	296	ASP	O-C-N	5.72	131.85	122.70
1	B	297	VAL	N-CA-C	-5.67	95.70	111.00
1	B	296	ASP	CA-C-N	-5.52	105.06	117.20
1	B	296	ASP	CB-CG-OD2	5.50	123.25	118.30
1	A	356	ASP	CB-CG-OD2	5.50	123.25	118.30
1	A	296	ASP	CB-CG-OD2	5.50	123.25	118.30
1	A	296	ASP	CA-C-N	-5.43	105.24	117.20
1	A	312	ASP	CB-CG-OD2	5.40	123.16	118.30
1	B	178	ASP	CB-CG-OD2	5.35	123.11	118.30
1	B	432	ASP	CB-CG-OD2	5.33	123.10	118.30
1	B	183	ALA	CA-C-N	-5.18	105.80	117.20
1	A	166	ASP	CB-CG-OD2	5.18	122.96	118.30
1	B	58	ASP	CB-CG-OD2	5.15	122.93	118.30
1	B	166	ASP	CB-CG-OD2	5.13	122.92	118.30
1	B	312	ASP	CB-CG-OD2	5.10	122.89	118.30
1	A	178	ASP	CB-CG-OD2	5.05	122.84	118.30
1	B	60	ARG	CA-C-N	-5.04	106.11	117.20

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	201	GLY	Peptide
1	A	296	ASP	Mainchain

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Mol	Chain	Res	Type	Group
1	A	298	GLU	Mainchain
1	B	201	GLY	Peptide
1	B	296	ASP	Mainchain
1	B	298	GLU	Mainchain

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	3334	0	3200	69	0
1	B	3334	0	3200	59	0
2	A	69	0	42	0	0
2	B	69	0	42	2	0
3	A	103	0	0	4	0
3	B	100	0	0	1	0
All	All	7009	0	6484	128	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 10.

All (128) 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:37:ASN:HD21	1:B:39:LYS:HZ3	1.13	0.92
1:B:37:ASN:HD21	1:B:39:LYS:NZ	1.70	0.88
1:A:37:ASN:HD21	1:A:39:LYS:NZ	1.73	0.86
1:A:37:ASN:HD21	1:A:39:LYS:HZ3	1.26	0.80
1:A:102:THR:CG2	1:A:104:LEU:H	1.95	0.78
1:A:102:THR:HG22	1:A:105:THR:H	1.50	0.77
1:B:37:ASN:H	1:B:37:ASN:HD22	1.31	0.77
1:A:194:ASP:HB3	1:A:196:ASN:H	1.50	0.76
1:B:232:THR:HG21	1:B:261:SER:H	1.50	0.75
1:A:232:THR:HG23	1:A:243:THR:OG1	1.86	0.75
1:B:232:THR:HG23	1:B:243:THR:OG1	1.87	0.74
1:B:102:THR:CG2	1:B:104:LEU:H	2.00	0.74
1:A:232:THR:CG2	1:A:243:THR:OG1	2.37	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:THR:HG21	1:A:261:SER:H	1.52	0.73
1:B:232:THR:CG2	1:B:243:THR:OG1	2.38	0.72
1:B:102:THR:HG22	1:B:105:THR:H	1.55	0.71
1:B:250:ASN:H	1:B:250:ASN:HD22	1.38	0.71
1:B:194:ASP:HB3	1:B:196:ASN:H	1.55	0.71
1:A:73:HIS:HD2	1:A:77:GLU:OE2	1.76	0.69
1:A:102:THR:HG22	1:A:104:LEU:H	1.56	0.68
1:A:250:ASN:H	1:A:250:ASN:HD22	1.41	0.68
1:B:77:GLU:OE1	1:B:102:THR:HG21	1.94	0.67
1:A:232:THR:HG21	1:A:260:ALA:HA	1.76	0.67
1:B:73:HIS:HD2	1:B:77:GLU:OE2	1.78	0.66
1:B:102:THR:HG22	1:B:104:LEU:H	1.60	0.66
1:A:151:HIS:HD2	1:A:320:GLU:OE1	1.79	0.66
1:B:151:HIS:HD2	1:B:320:GLU:OE1	1.79	0.65
1:B:232:THR:HG21	1:B:260:ALA:HA	1.79	0.65
1:B:365:ILE:HG23	1:B:370:GLU:HG2	1.77	0.64
1:A:77:GLU:OE1	1:A:102:THR:HG21	1.97	0.64
1:A:181:ILE:HD11	1:A:247:PHE:CE1	2.33	0.64
1:A:365:ILE:HG23	1:A:370:GLU:HG2	1.80	0.63
1:A:94:VAL:HG13	3:A:456:HOH:O	1.98	0.62
1:B:178:ASP:HB2	1:B:193:LYS:HB2	1.82	0.62
1:A:37:ASN:H	1:A:37:ASN:HD22	1.46	0.61
1:B:224:GLU:HA	1:B:250:ASN:HD21	1.65	0.61
1:B:225:HIS:H	1:B:250:ASN:ND2	1.99	0.61
1:A:225:HIS:H	1:A:250:ASN:ND2	1.98	0.60
1:A:102:THR:HG23	1:A:104:LEU:H	1.64	0.60
1:B:94:VAL:HG13	3:B:457:HOH:O	2.01	0.60
1:B:74:TRP:CD2	1:B:75:PRO:HA	2.37	0.60
1:B:6:ILE:HD13	1:B:292:VAL:HG13	1.84	0.60
1:A:182:TYR:HA	1:A:213:VAL:O	2.02	0.59
1:B:58:ASP:OD2	1:B:63:GLU:HG2	2.02	0.59
1:A:224:GLU:HA	1:A:250:ASN:HD21	1.67	0.58
1:B:181:ILE:HD11	1:B:247:PHE:CE1	2.38	0.58
1:B:37:ASN:H	1:B:37:ASN:ND2	1.99	0.58
1:A:74:TRP:CD2	1:A:75:PRO:HA	2.39	0.58
1:B:432:ASP:N	2:B:455:XMM:O6	2.36	0.58
1:B:61:GLU:C	1:B:63:GLU:H	2.07	0.57
1:B:102:THR:HG23	1:B:104:LEU:H	1.69	0.56
1:A:175:TRP:CE2	1:A:195:ALA:HB2	2.40	0.56
1:B:6:ILE:CD1	1:B:292:VAL:HG13	2.35	0.56
1:A:178:ASP:HB2	1:A:193:LYS:HB2	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:250:ASN:H	1:A:250:ASN:ND2	2.04	0.55
1:A:58:ASP:OD2	1:A:63:GLU:HG2	2.06	0.55
1:B:37:ASN:N	1:B:37:ASN:HD22	2.03	0.55
1:B:250:ASN:H	1:B:250:ASN:ND2	2.03	0.54
1:B:232:THR:HG21	1:B:261:SER:N	2.21	0.54
1:B:232:THR:HB	1:B:244:SER:OG	2.08	0.54
1:A:33:HIS:HD2	3:A:488:HOH:O	1.91	0.52
1:A:37:ASN:N	1:A:37:ASN:HD22	2.05	0.52
1:B:225:HIS:H	1:B:250:ASN:HD21	1.56	0.52
1:A:185:ALA:HA	1:A:259:ILE:HG13	1.91	0.52
1:B:73:HIS:CD2	1:B:77:GLU:OE2	2.61	0.52
1:A:73:HIS:CD2	1:A:77:GLU:OE2	2.59	0.52
1:A:232:THR:HB	1:A:244:SER:OG	2.10	0.52
1:B:224:GLU:HG3	1:B:251:VAL:HB	1.92	0.52
1:A:225:HIS:H	1:A:250:ASN:HD21	1.56	0.52
1:A:224:GLU:HG3	1:A:251:VAL:HB	1.92	0.50
1:B:232:THR:HG23	1:B:261:SER:O	2.12	0.49
1:B:232:THR:CG2	1:B:261:SER:H	2.23	0.49
1:A:37:ASN:H	1:A:37:ASN:ND2	2.11	0.49
1:B:37:ASN:ND2	1:B:39:LYS:NZ	2.50	0.49
1:A:37:ASN:ND2	1:A:39:LYS:NZ	2.54	0.48
1:A:360:LYS:HE2	3:A:484:HOH:O	2.14	0.48
1:A:185:ALA:HA	1:A:259:ILE:CD1	2.45	0.47
1:B:177:LYS:HE2	1:B:178:ASP:OD2	2.15	0.47
1:A:185:ALA:HA	1:A:259:ILE:HD12	1.97	0.47
1:A:393:SER:HA	1:A:406:PHE:O	2.14	0.47
1:A:185:ALA:HA	1:A:259:ILE:CG1	2.45	0.47
1:B:433:TYR:H	2:B:455:XMM:HO6	1.61	0.47
1:A:232:THR:HG21	1:A:261:SER:N	2.24	0.46
1:B:424:VAL:HG11	1:B:442:LYS:HD2	1.97	0.46
1:B:244:SER:HA	1:B:257:PHE:O	2.16	0.46
1:A:232:THR:HG23	1:A:261:SER:O	2.15	0.46
1:A:25:HIS:HD2	1:A:45:ASP:OD2	1.99	0.46
1:A:45:ASP:OD2	1:A:49:ASP:HB2	2.16	0.46
1:A:226:LEU:HB3	1:A:274:VAL:HB	1.97	0.45
1:B:25:HIS:HD2	1:B:45:ASP:OD2	1.99	0.45
1:A:225:HIS:N	1:A:250:ASN:HD21	2.15	0.45
1:A:424:VAL:HG11	1:A:442:LYS:HD2	1.98	0.45
1:A:244:SER:HA	1:A:257:PHE:O	2.15	0.45
1:B:393:SER:HA	1:B:406:PHE:O	2.16	0.45
1:A:267:ILE:HD12	1:A:268:PRO:HD2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:225:HIS:N	1:B:250:ASN:HD21	2.15	0.44
1:B:29:GLU:OE1	1:B:71:LYS:HE3	2.16	0.44
1:A:102:THR:HG22	1:A:104:LEU:N	2.29	0.44
1:B:61:GLU:C	1:B:63:GLU:N	2.71	0.44
1:A:443:PRO:O	1:A:444:ASP:HB2	2.18	0.44
1:A:278:HIS:HD2	3:A:495:HOH:O	1.99	0.44
1:A:232:THR:CG2	1:A:261:SER:H	2.25	0.44
1:A:324:GLN:HB3	1:A:339:LYS:HB2	1.99	0.43
1:A:28:THR:HA	1:A:72:ILE:HB	2.00	0.43
1:B:324:GLN:HB3	1:B:339:LYS:HB2	2.00	0.43
1:B:28:THR:HA	1:B:72:ILE:HB	2.01	0.43
1:B:181:ILE:HD11	1:B:247:PHE:CZ	2.54	0.42
1:A:70:ILE:N	1:A:70:ILE:HD12	2.33	0.42
1:B:267:ILE:HD12	1:B:268:PRO:HD2	2.00	0.42
1:A:316:TYR:CD1	1:A:346:SER:HB3	2.55	0.42
1:A:94:VAL:HG22	1:A:140:ASP:HA	2.02	0.42
1:A:37:ASN:ND2	1:A:37:ASN:N	2.67	0.42
1:A:160:GLY:HA2	1:A:233:TYR:CG	2.55	0.42
1:B:316:TYR:CD1	1:B:346:SER:HB3	2.54	0.42
1:B:443:PRO:O	1:B:444:ASP:HB2	2.20	0.42
1:A:102:THR:HG22	1:A:105:THR:N	2.27	0.41
1:B:226:LEU:HB3	1:B:274:VAL:HB	2.01	0.41
1:A:80:LYS:HE2	1:A:101:ILE:HG22	2.03	0.41
1:A:385:LYS:HZ2	1:A:385:LYS:HG2	1.66	0.41
1:B:318:ALA:HB2	1:B:322:ILE:HG13	2.02	0.41
1:B:252:THR:OG1	1:B:253:LYS:N	2.53	0.41
1:A:7:SER:O	1:A:269:ILE:HA	2.20	0.41
1:A:29:GLU:OE1	1:A:71:LYS:HE3	2.20	0.41
1:B:15:GLY:O	1:B:134:ARG:HD3	2.21	0.41
1:A:149:VAL:HB	1:A:150:PRO:HD2	2.03	0.40
1:B:7:SER:O	1:B:269:ILE:HA	2.21	0.40
1:A:15:GLY:O	1:A:134:ARG:HD3	2.21	0.40
1:A:124:ALA:HB3	1:A:155:SER:HB3	2.02	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	440/447 (98%)	416 (94%)	22 (5%)	2 (0%)	34	55
1	B	440/447 (98%)	418 (95%)	18 (4%)	4 (1%)	21	37
All	All	880/894 (98%)	834 (95%)	40 (4%)	6 (1%)	26	46

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	298	GLU
1	B	298	GLU
1	B	4	GLY
1	B	62	ASN
1	B	194	ASP
1	A	4	GLY

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	353/358 (99%)	326 (92%)	27 (8%)	16	30
1	B	353/358 (99%)	325 (92%)	28 (8%)	15	28
All	All	706/716 (99%)	651 (92%)	55 (8%)	16	29

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	62	ASN
1	A	68	LYS
1	A	88	ASP
1	A	93	LEU
1	A	94	VAL
1	A	102	THR
1	A	114	SER
1	A	117	GLU
1	A	134	ARG
1	A	194	ASP
1	A	215	LYS
1	A	226	LEU
1	A	232	THR
1	A	250	ASN
1	A	285	LEU
1	A	293	LYS
1	A	296	ASP
1	A	303	ILE
1	A	323	ASN
1	A	336	VAL
1	A	340	ASP
1	A	373	THR
1	A	385	LYS
1	A	387	VAL
1	A	416	ILE
1	A	444	ASP
1	B	3	LYS
1	B	37	ASN
1	B	59	PRO
1	B	62	ASN
1	B	68	LYS
1	B	76	THR
1	B	88	ASP
1	B	93	LEU
1	B	94	VAL
1	B	102	THR
1	B	117	GLU
1	B	134	ARG
1	B	215	LYS
1	B	226	LEU
1	B	232	THR
1	B	250	ASN

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Mol	Chain	Res	Type
1	B	285	LEU
1	B	292	VAL
1	B	293	LYS
1	B	296	ASP
1	B	303	ILE
1	B	323	ASN
1	B	336	VAL
1	B	340	ASP
1	B	373	THR
1	B	385	LYS
1	B	387	VAL
1	B	416	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	25	HIS
1	A	33	HIS
1	A	37	ASN
1	A	73	HIS
1	A	151	HIS
1	A	188	ASN
1	A	235	ASN
1	A	250	ASN
1	A	278	HIS
1	A	323	ASN
1	B	24	ASN
1	B	25	HIS
1	B	33	HIS
1	B	37	ASN
1	B	62	ASN
1	B	73	HIS
1	B	151	HIS
1	B	188	ASN
1	B	235	ASN
1	B	250	ASN
1	B	323	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

6 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	XMM	A	450	-	24,25,25	5.14	8 (33%)	27,37,37	9.25	14 (51%)
2	XMM	A	451	-	24,25,25	5.07	8 (33%)	27,37,37	8.91	13 (48%)
2	XMM	A	452	-	24,25,25	4.80	8 (33%)	27,37,37	8.38	12 (44%)
2	XMM	B	453	-	24,25,25	4.81	9 (37%)	27,37,37	8.78	13 (48%)
2	XMM	B	454	-	24,25,25	5.04	8 (33%)	27,37,37	8.94	9 (33%)
2	XMM	B	455	-	24,25,25	5.18	8 (33%)	27,37,37	9.20	12 (44%)

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	XMM	A	450	-	-	0/4/26/26	0/3/3/3
2	XMM	A	451	-	-	0/4/26/26	0/3/3/3
2	XMM	A	452	-	-	0/4/26/26	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	XMM	B	453	-	-	0/4/26/26	0/3/3/3
2	XMM	B	454	-	-	0/4/26/26	0/3/3/3
2	XMM	B	455	-	-	0/4/26/26	0/3/3/3

All (49) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	455	XMM	BR-C13	-18.02	1.45	1.89
2	A	450	XMM	BR-C13	-17.89	1.45	1.89
2	A	451	XMM	BR-C13	-17.64	1.46	1.89
2	B	454	XMM	BR-C13	-17.42	1.46	1.89
2	B	453	XMM	BR-C13	-16.66	1.48	1.89
2	A	452	XMM	BR-C13	-16.42	1.49	1.89
2	B	453	XMM	C10-CL	-6.88	1.46	1.73
2	B	454	XMM	C10-CL	-6.62	1.47	1.73
2	A	452	XMM	C10-CL	-6.60	1.47	1.73
2	A	451	XMM	C10-CL	-6.28	1.48	1.73
2	A	450	XMM	C10-CL	-6.18	1.49	1.73
2	B	455	XMM	C10-CL	-6.10	1.49	1.73
2	B	455	XMM	C14-C13	-5.21	1.27	1.39
2	A	450	XMM	C14-C13	-5.10	1.27	1.39
2	B	454	XMM	C14-C13	-4.68	1.28	1.39
2	A	451	XMM	C14-C13	-4.51	1.28	1.39
2	B	454	XMM	C11-N1	-4.48	1.27	1.36
2	B	453	XMM	C14-C13	-4.46	1.29	1.39
2	A	451	XMM	C11-N1	-4.43	1.27	1.36
2	B	453	XMM	C11-N1	-4.32	1.27	1.36
2	A	452	XMM	C11-N1	-4.28	1.27	1.36
2	A	452	XMM	C14-C13	-4.18	1.29	1.39
2	B	455	XMM	C11-N1	-3.86	1.28	1.36
2	A	450	XMM	C11-N1	-3.58	1.29	1.36
2	B	453	XMM	C12-C14	-3.56	1.28	1.36
2	A	452	XMM	C12-C14	-3.28	1.29	1.36
2	B	455	XMM	C11-C7	-3.01	1.29	1.37
2	B	454	XMM	C11-C7	-3.00	1.29	1.37
2	A	450	XMM	C11-C7	-2.87	1.30	1.37
2	A	451	XMM	C11-C7	-2.83	1.30	1.37
2	B	453	XMM	C11-C7	-2.82	1.30	1.37
2	B	454	XMM	C12-C14	-2.81	1.30	1.36
2	B	455	XMM	C12-C14	-2.80	1.30	1.36
2	A	450	XMM	C12-C14	-2.70	1.30	1.36
2	A	451	XMM	C12-C14	-2.69	1.30	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	452	XMM	C11-C7	-2.60	1.30	1.37
2	B	453	XMM	O6-C6	2.20	1.51	1.42
2	B	453	XMM	C7-C8	8.51	1.51	1.40
2	A	452	XMM	C7-C8	8.98	1.51	1.40
2	B	454	XMM	C7-C8	9.04	1.52	1.40
2	A	451	XMM	C7-C8	9.17	1.52	1.40
2	B	453	XMM	C10-C13	9.39	1.53	1.40
2	A	452	XMM	C10-C13	9.80	1.54	1.40
2	B	455	XMM	C7-C8	10.21	1.53	1.40
2	A	450	XMM	C7-C8	10.34	1.53	1.40
2	A	450	XMM	C10-C13	10.38	1.54	1.40
2	B	455	XMM	C10-C13	10.55	1.55	1.40
2	B	454	XMM	C10-C13	10.64	1.55	1.40
2	A	451	XMM	C10-C13	10.96	1.55	1.40

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	454	XMM	C14-C13-C10	-19.16	100.77	120.23
2	A	451	XMM	C14-C13-C10	-19.15	100.78	120.23
2	B	453	XMM	C14-C13-C10	-18.02	101.92	120.23
2	A	450	XMM	C14-C13-C10	-17.73	102.22	120.23
2	A	452	XMM	C14-C13-C10	-17.43	102.52	120.23
2	B	455	XMM	C14-C13-C10	-17.42	102.53	120.23
2	B	453	XMM	C14-C12-C9	-12.69	107.10	120.88
2	A	450	XMM	C14-C12-C9	-12.52	107.29	120.88
2	B	455	XMM	C14-C12-C9	-12.38	107.44	120.88
2	A	452	XMM	C14-C12-C9	-12.12	107.72	120.88
2	B	454	XMM	C14-C12-C9	-11.33	108.58	120.88
2	A	451	XMM	C14-C12-C9	-11.15	108.77	120.88
2	A	450	XMM	C6-C5-C4	-5.70	98.96	113.02
2	B	455	XMM	C6-C5-C4	-4.10	102.89	113.02
2	B	455	XMM	C13-C10-C8	-4.10	116.94	121.28
2	A	450	XMM	C13-C10-C8	-3.96	117.09	121.28
2	B	454	XMM	C12-C9-C8	-3.67	114.81	121.11
2	A	451	XMM	C12-C9-C8	-3.57	114.99	121.11
2	B	455	XMM	O1-C1-C2	-3.32	101.73	107.12
2	A	451	XMM	C13-C10-C8	-3.29	117.80	121.28
2	B	453	XMM	O3-C3-C4	-3.15	103.25	110.34
2	B	453	XMM	C13-C10-C8	-3.08	118.02	121.28
2	B	454	XMM	C13-C10-C8	-3.04	118.06	121.28
2	A	451	XMM	O6-C6-C5	-2.95	101.59	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	450	XMM	C12-C9-C8	-2.87	116.19	121.11
2	B	453	XMM	O5-C1-O1	-2.85	100.78	108.39
2	A	451	XMM	C6-C5-C4	-2.82	106.05	113.02
2	A	452	XMM	C13-C10-C8	-2.82	118.30	121.28
2	B	453	XMM	C12-C9-C8	-2.81	116.29	121.11
2	A	452	XMM	C12-C9-C8	-2.78	116.35	121.11
2	B	455	XMM	C12-C9-C8	-2.74	116.41	121.11
2	A	452	XMM	O5-C1-O1	-2.70	101.18	108.39
2	B	454	XMM	O4-C4-C3	-2.51	104.69	110.34
2	B	455	XMM	O2-C2-C3	-2.50	104.71	110.34
2	A	452	XMM	O3-C3-C4	-2.48	104.75	110.34
2	A	450	XMM	O1-C1-C2	-2.39	103.23	107.12
2	A	450	XMM	O2-C2-C3	-2.34	105.07	110.34
2	A	450	XMM	C4-C3-C2	-2.22	106.64	110.79
2	B	453	XMM	O5-C5-C4	-2.15	105.65	109.68
2	A	451	XMM	O3-C3-C4	-2.03	105.76	110.34
2	A	452	XMM	O5-C5-C6	2.05	111.52	106.36
2	B	454	XMM	O3-C3-C2	2.06	114.98	110.34
2	B	453	XMM	O1-C1-C2	2.31	110.89	107.12
2	A	451	XMM	O5-C5-C4	2.39	114.17	109.68
2	B	455	XMM	O5-C5-C6	2.44	112.52	106.36
2	A	451	XMM	O3-C3-C2	2.58	116.14	110.34
2	A	452	XMM	O1-C1-C2	2.60	111.35	107.12
2	A	450	XMM	O5-C5-C4	2.62	114.59	109.68
2	A	450	XMM	C1-O5-C5	2.79	119.16	113.75
2	A	451	XMM	C7-O1-C1	2.80	123.50	118.24
2	B	453	XMM	O5-C5-C6	2.91	113.72	106.36
2	B	455	XMM	BR-C13-C14	3.38	124.55	117.81
2	A	450	XMM	BR-C13-C14	3.61	125.01	117.81
2	B	454	XMM	BR-C13-C14	4.76	127.29	117.81
2	A	451	XMM	BR-C13-C14	4.97	127.70	117.81
2	B	453	XMM	BR-C13-C14	5.37	128.51	117.81
2	A	452	XMM	BR-C13-C14	5.79	129.34	117.81
2	B	453	XMM	C7-O1-C1	7.08	131.54	118.24
2	A	452	XMM	C7-O1-C1	7.77	132.83	118.24
2	A	450	XMM	C7-O1-C1	9.55	136.17	118.24
2	B	455	XMM	C7-O1-C1	10.28	137.54	118.24
2	A	452	XMM	BR-C13-C10	10.49	128.14	121.12
2	B	453	XMM	BR-C13-C10	12.63	129.56	121.12
2	A	451	XMM	BR-C13-C10	15.52	131.49	121.12
2	B	454	XMM	BR-C13-C10	16.18	131.94	121.12
2	A	450	XMM	BR-C13-C10	17.34	132.71	121.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	455	XMM	BR-C13-C10	17.61	132.89	121.12
2	A	452	XMM	C12-C14-C13	34.57	151.79	119.63
2	B	453	XMM	C12-C14-C13	35.97	153.10	119.63
2	A	451	XMM	C12-C14-C13	36.14	153.26	119.63
2	B	455	XMM	C12-C14-C13	36.37	153.47	119.63
2	B	454	XMM	C12-C14-C13	36.49	153.59	119.63
2	A	450	XMM	C12-C14-C13	36.62	153.70	119.63

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	455	XMM	2	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 [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	439/447 (98%)	-0.25	3 (0%) 89 90	9, 24, 41, 54	4 (0%)
1	B	439/447 (98%)	-0.26	3 (0%) 89 90	9, 24, 41, 54	2 (0%)
All	All	878/894 (98%)	-0.26	6 (0%) 89 90	9, 24, 41, 54	6 (0%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	222	ASN	3.0
1	A	222	ASN	2.4
1	A	183	ALA	2.3
1	B	183	ALA	2.2
1	B	35	LYS	2.2
1	A	185	ALA	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	XMM	A	451	23/23	0.88	0.17	1.14	8,18,26,36	0
2	XMM	B	455	23/23	0.87	0.16	1.02	11,22,30,35	0
2	XMM	A	450	23/23	0.90	0.15	0.76	10,21,32,38	0
2	XMM	A	452	23/23	0.83	0.17	0.73	11,20,25,29	0
2	XMM	B	453	23/23	0.83	0.17	0.65	12,21,27,32	0
2	XMM	B	454	23/23	0.87	0.16	0.64	9,18,24,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.