



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:17 PM GMT

PDB ID : 4ZTL
Title : Irak4-inhibitor co-structure
Authors : Fischmann, T.O.
Deposited on : 2015-05-14
Resolution : 2.39 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

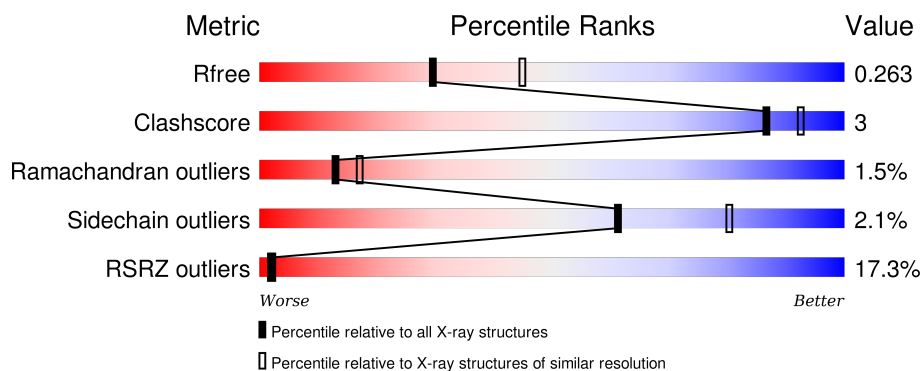
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (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 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	301	<div> <div>21%</div> <div>87%</div> <div>6% • 6%</div> </div>
1	B	301	<div> <div>19%</div> <div>84%</div> <div>6% 10%</div> </div>
1	C	301	<div> <div>14%</div> <div>84%</div> <div>10% • 5%</div> </div>
1	D	301	<div> <div>11%</div> <div>86%</div> <div>9% •</div> </div>

2 Entry composition [i](#)

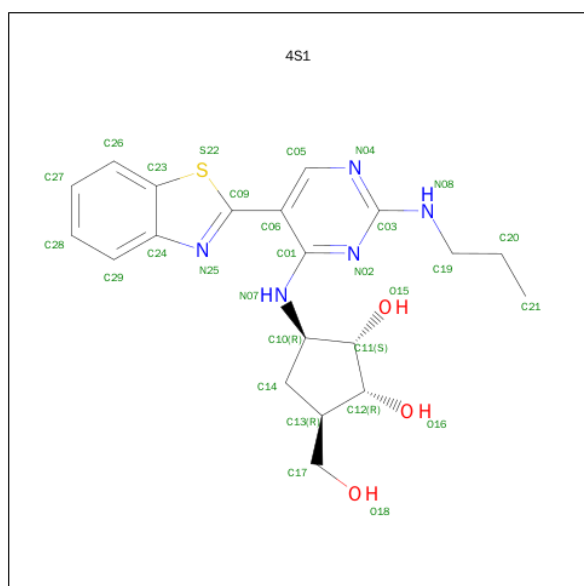
There are 3 unique types of molecules in this entry. The entry contains 9076 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Interleukin-1 receptor-associated kinase 4.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	283	Total	C	N	O	P	S	0	1	3
			2230	1398	372	442	3	15			
1	B	272	Total	C	N	O	P	S	0	1	4
			2124	1332	355	419	3	15			
1	C	285	Total	C	N	O	P	S	0	0	2
			2239	1405	378	439	3	14			
1	D	288	Total	C	N	O	P	S	0	0	4
			2259	1414	382	446	3	14			

- Molecule 2 is (1R,2S,3R,5R)-3-{{5-(1,3-benzothiazol-2-yl)-2-(propylamino)pyrimidin-4-yl}amino}-5-(hydroxymethyl)cyclopentane-1,2-diol (three-letter code: 4S1) (formula: C₂₀H₂₅N₅O₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			29	20	5	3	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	S	0	0
			29	20	5	3	1		
2	C	1	Total	C	N	O	S	0	0
			29	20	5	3	1		
2	D	1	Total	C	N	O	S	0	0
			29	20	5	3	1		

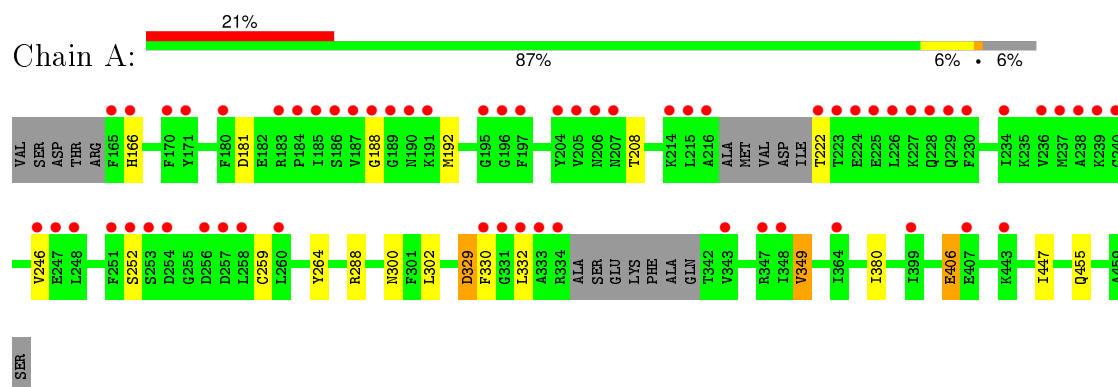
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	19	Total	O	0	0
			19	19		
3	B	19	Total	O	0	0
			19	19		
3	C	35	Total	O	0	0
			35	35		
3	D	35	Total	O	0	0
			35	35		

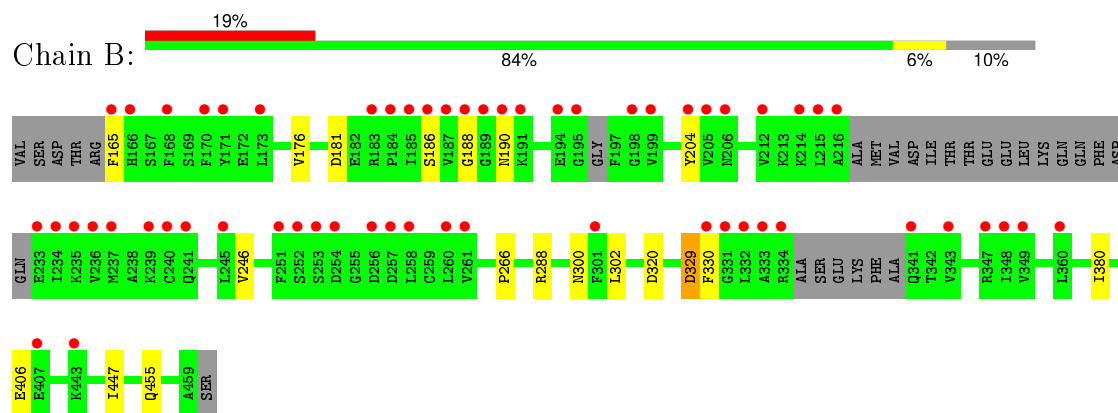
3 Residue-property plots [i](#)

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.

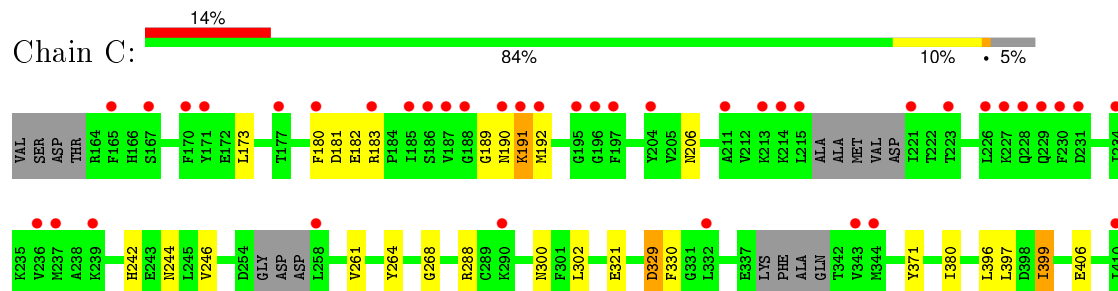
- Molecule 1: Interleukin-1 receptor-associated kinase 4

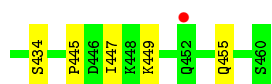


- Molecule 1: Interleukin-1 receptor-associated kinase 4

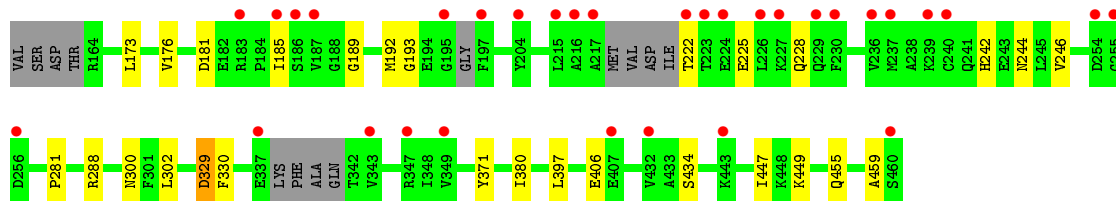
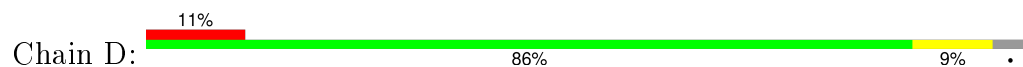


- Molecule 1: Interleukin-1 receptor-associated kinase 4





- Molecule 1: Interleukin-1 receptor-associated kinase 4



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	145.18Å 139.39Å 87.98Å 90.00° 125.93° 90.00°	Depositor
Resolution (Å)	28.31 – 2.39 28.31 – 2.39	Depositor EDS
% Data completeness (in resolution range)	98.0 (28.31-2.39) 92.4 (28.31-2.39)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 2.39Å)	Xtriage
Refinement program	BUSTER	Depositor
R, R_{free}	0.225 , 0.251 0.237 , 0.263	Depositor DCC
R_{free} test set	2817 reflections (5.41%)	DCC
Wilson B-factor (Å ²)	48.7	Xtriage
Anisotropy	0.467	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.0	EDS
Estimated twinning fraction	0.317 for -h-2*l,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 54882 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9076	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.67 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 5.4411e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: TPO, 4S1, SEP

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	0.48	0/2234	0.66	0/3010
1	B	0.48	0/2124	0.66	0/2860
1	C	0.52	0/2242	0.69	0/3019
1	D	0.52	0/2262	0.67	0/3047
All	All	0.50	0/8862	0.67	0/11936

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2230	0	2186	9	0
1	B	2124	0	2084	8	0
1	C	2239	0	2207	18	0
1	D	2259	0	2218	14	0
2	A	29	0	25	2	0
2	B	29	0	25	2	0
2	C	29	0	25	3	0
2	D	29	0	25	3	0
3	A	19	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	19	0	0	0	0
3	C	35	0	0	0	0
3	D	35	0	0	0	0
All	All	9076	0	8795	48	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 3.

All (48) 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:D:242:HIS:CD2	1:D:244:ASN:H	2.17	0.62
1:C:242:HIS:CD2	1:C:244:ASN:H	2.21	0.58
1:B:181:ASP:HB3	1:B:190:ASN:HB2	1.87	0.57
1:C:246:VAL:HG21	2:C:501:4S1:H22	1.87	0.57
1:D:242:HIS:HD2	1:D:244:ASN:H	1.54	0.56
1:B:246:VAL:HG21	2:B:501:4S1:H22	1.87	0.56
1:C:182:GLU:HA	1:C:191:LYS:HB2	1.92	0.51
1:D:222:THR:HB	1:D:225:GLU:HB2	1.92	0.51
1:D:246:VAL:HG21	2:D:501:4S1:H22	1.92	0.51
1:D:288:ARG:HB3	1:D:380:ILE:HG23	1.93	0.50
1:A:332:LEU:HD13	1:A:349:VAL:HG13	1.93	0.50
1:C:268:GLY:HA2	2:C:501:4S1:H16	1.93	0.49
1:C:288:ARG:HB3	1:C:380:ILE:HG23	1.94	0.49
1:C:246:VAL:HG21	2:C:501:4S1:C26	2.42	0.49
1:A:302:LEU:HD11	1:A:330:PHE:HE1	1.78	0.48
1:D:193:GLY:HA3	2:D:501:4S1:H8	1.95	0.48
1:C:242:HIS:HD2	1:C:244:ASN:H	1.60	0.48
1:D:300:ASN:HA	1:D:447:ILE:HG21	1.97	0.47
1:A:246:VAL:HG21	2:A:501:4S1:C26	2.45	0.47
1:C:181:ASP:HB3	1:C:190:ASN:HB2	1.96	0.46
1:B:302:LEU:HD11	1:B:330:PHE:HE1	1.81	0.46
1:A:192:MET:CE	1:A:264:TYR:HE1	2.29	0.46
1:B:246:VAL:HG21	2:B:501:4S1:C26	2.46	0.46
1:B:266:PRO:HD2	1:B:320:ASP:HA	1.98	0.45
1:C:300:ASN:HA	1:C:447:ILE:HG21	1.98	0.45
1:C:445:PRO:HB3	1:C:449:LYS:HD3	1.97	0.45
1:A:300:ASN:HA	1:A:447:ILE:HG21	1.99	0.45
1:C:321:GLU:HG3	1:D:281:PRO:HD3	1.98	0.45
1:C:183:ARG:HB3	1:C:189:GLY:HA3	1.99	0.45
1:C:302:LEU:HD11	1:C:330:PHE:HE1	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:302:LEU:HD11	1:D:330:PHE:HE1	1.83	0.44
1:C:396:LEU:HD12	1:C:399:ILE:HD13	1.99	0.44
1:A:246:VAL:HG21	2:A:501:4S1:H22	1.99	0.44
1:B:300:ASN:HA	1:B:447:ILE:HG21	1.99	0.44
1:D:185:ILE:HD11	1:D:192:MET:HG2	2.01	0.42
1:A:332:LEU:HD22	1:A:349:VAL:HG11	2.02	0.42
1:C:371:TYR:OH	1:C:397:LEU:HD22	2.20	0.42
1:B:176:VAL:HB	1:B:204:TYR:H	1.84	0.42
1:D:246:VAL:HG21	2:D:501:4S1:C26	2.50	0.41
1:C:173:LEU:HD21	1:C:261:VAL:HG21	2.02	0.41
1:C:192:MET:CE	1:C:264:TYR:HE1	2.33	0.41
1:C:396:LEU:O	1:C:399:ILE:HB	2.20	0.41
1:B:288:ARG:HB3	1:B:380:ILE:HG23	2.03	0.41
1:D:185:ILE:HA	1:D:189:GLY:O	2.21	0.41
1:D:173:LEU:HA	1:D:176:VAL:HG22	2.02	0.41
1:A:288:ARG:HB3	1:A:380:ILE:HG23	2.02	0.41
1:D:371:TYR:OH	1:D:397:LEU:HD22	2.21	0.40
1:A:252:SER:HB3	1:A:259:CYS:HB2	2.04	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	276/301 (92%)	266 (96%)	7 (2%)	3 (1%)	17	25
1	B	262/301 (87%)	247 (94%)	11 (4%)	4 (2%)	13	17
1	C	275/301 (91%)	260 (94%)	10 (4%)	5 (2%)	11	13
1	D	278/301 (92%)	265 (95%)	9 (3%)	4 (1%)	14	19
All	All	1091/1204 (91%)	1038 (95%)	37 (3%)	16 (2%)	13	17

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	406	GLU
1	B	406	GLU
1	C	406	GLU
1	D	406	GLU
1	D	459	ALA
1	D	181	ASP
1	A	329	ASP
1	B	329	ASP
1	C	191	LYS
1	C	329	ASP
1	D	329	ASP
1	A	188	GLY
1	B	188	GLY
1	C	206	ASN
1	C	180	PHE
1	B	186	SER

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	244/259 (94%)	236 (97%)	8 (3%)	45	66
1	B	232/259 (90%)	229 (99%)	3 (1%)	76	89
1	C	244/259 (94%)	240 (98%)	4 (2%)	70	86
1	D	246/259 (95%)	241 (98%)	5 (2%)	63	81
All	All	966/1036 (93%)	946 (98%)	20 (2%)	61	80

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	166	HIS
1	A	181	ASP
1	A	208	THR
1	A	222	THR
1	A	329	ASP
1	A	349	VAL

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Mol	Chain	Res	Type
1	A	406	GLU
1	A	455	GLN
1	B	165	PHE
1	B	329	ASP
1	B	455	GLN
1	C	329	ASP
1	C	399	ILE
1	C	434	SER
1	C	455	GLN
1	D	228	GLN
1	D	329	ASP
1	D	434	SER
1	D	449	LYS
1	D	455	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	394	GLN
1	B	242	HIS
1	B	394	GLN
1	C	242	HIS
1	C	394	GLN
1	D	228	GLN
1	D	242	HIS
1	D	394	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

12 non-standard protein/DNA/RNA residues 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$
1	TPO	A	342	1	7,10,11	0.81	0	10,14,16	1.20	1 (10%)
1	TPO	A	345	1	7,10,11	0.91	0	10,14,16	1.39	1 (10%)
1	SEP	A	346	1	7,9,10	0.98	0	8,12,14	1.17	1 (12%)
1	TPO	B	342	1	7,10,11	0.78	0	10,14,16	1.21	1 (10%)
1	TPO	B	345	1	7,10,11	0.86	0	10,14,16	1.39	1 (10%)
1	SEP	B	346	1	7,9,10	0.89	0	8,12,14	1.14	1 (12%)
1	TPO	C	342	1	7,10,11	0.87	0	10,14,16	1.17	0
1	TPO	C	345	1	7,10,11	1.34	1 (14%)	10,14,16	1.54	1 (10%)
1	SEP	C	346	1	7,9,10	0.90	0	8,12,14	1.43	2 (25%)
1	TPO	D	342	1	7,10,11	0.78	0	10,14,16	1.18	0
1	TPO	D	345	1	7,10,11	0.81	0	10,14,16	1.41	1 (10%)
1	SEP	D	346	1	7,9,10	1.07	0	8,12,14	1.09	1 (12%)

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
1	TPO	A	342	1	-	0/8/11/13	0/0/0/0
1	TPO	A	345	1	-	1/8/11/13	0/0/0/0
1	SEP	A	346	1	-	0/5/8/10	0/0/0/0
1	TPO	B	342	1	-	0/8/11/13	0/0/0/0
1	TPO	B	345	1	-	1/8/11/13	0/0/0/0
1	SEP	B	346	1	-	0/5/8/10	0/0/0/0
1	TPO	C	342	1	-	0/8/11/13	0/0/0/0
1	TPO	C	345	1	-	1/8/11/13	0/0/0/0
1	SEP	C	346	1	-	0/5/8/10	0/0/0/0
1	TPO	D	342	1	-	0/8/11/13	0/0/0/0
1	TPO	D	345	1	-	1/8/11/13	0/0/0/0
1	SEP	D	346	1	-	0/5/8/10	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	345	TPO	CB-CA	2.77	1.59	1.54

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	346	SEP	O-C-CA	-2.46	119.12	125.72
1	D	346	SEP	O-C-CA	-2.25	119.69	125.72
1	B	346	SEP	O-C-CA	-2.25	119.69	125.72
1	C	346	SEP	O-C-CA	-2.22	119.77	125.72
1	B	342	TPO	P-OG1-CB	-2.12	112.16	121.42
1	A	342	TPO	P-OG1-CB	-2.07	112.35	121.42
1	B	345	TPO	O3P-P-OG1	2.41	113.83	106.62
1	A	345	TPO	O3P-P-OG1	2.42	113.87	106.62
1	D	345	TPO	O3P-P-OG1	2.59	114.37	106.62
1	C	346	SEP	OG-CB-CA	3.02	110.89	108.26
1	C	345	TPO	OG1-P-O1P	3.51	115.85	107.48

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	345	TPO	OG1-CB-CA-N
1	B	345	TPO	OG1-CB-CA-N
1	D	345	TPO	OG1-CB-CA-N
1	C	345	TPO	OG1-CB-CA-N

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates 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	4S1	A	501	-	29,32,32	1.87	11 (37%)	31,45,45	1.60	5 (16%)
2	4S1	B	501	-	29,32,32	1.96	11 (37%)	31,45,45	1.77	9 (29%)
2	4S1	C	501	-	29,32,32	1.89	9 (31%)	31,45,45	1.59	7 (22%)
2	4S1	D	501	-	29,32,32	1.79	9 (31%)	31,45,45	1.99	9 (29%)

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	4S1	A	501	-	-	0/10/30/30	0/4/4/4
2	4S1	B	501	-	-	0/10/30/30	0/4/4/4
2	4S1	C	501	-	-	0/10/30/30	0/4/4/4
2	4S1	D	501	-	-	0/10/30/30	0/4/4/4

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	4S1	C28-C27	2.05	1.43	1.38
2	C	501	4S1	C28-C27	2.07	1.43	1.38
2	A	501	4S1	C10-N07	2.12	1.48	1.45
2	B	501	4S1	C01-N02	2.20	1.38	1.34
2	A	501	4S1	C01-N02	2.25	1.38	1.34
2	A	501	4S1	C28-C27	2.25	1.43	1.38
2	D	501	4S1	C10-N07	2.25	1.48	1.45
2	D	501	4S1	C28-C29	2.31	1.41	1.36
2	B	501	4S1	C06-C01	2.38	1.47	1.42
2	A	501	4S1	C01-N07	2.38	1.38	1.35
2	D	501	4S1	C06-C01	2.40	1.47	1.42
2	A	501	4S1	C17-C13	2.40	1.57	1.52
2	B	501	4S1	C10-N07	2.41	1.48	1.45
2	D	501	4S1	C27-C26	2.46	1.42	1.36
2	A	501	4S1	C09-S22	2.49	1.77	1.73
2	C	501	4S1	C27-C26	2.50	1.42	1.36
2	B	501	4S1	C28-C27	2.52	1.44	1.38
2	C	501	4S1	C05-C06	2.53	1.43	1.39
2	C	501	4S1	C28-C29	2.54	1.42	1.36
2	C	501	4S1	C03-N04	2.56	1.37	1.34
2	B	501	4S1	C05-N04	2.59	1.39	1.34
2	B	501	4S1	C27-C26	2.70	1.42	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	4S1	C28-C29	2.79	1.43	1.36
2	B	501	4S1	C03-N08	2.82	1.38	1.34
2	A	501	4S1	C09-N25	2.90	1.35	1.31
2	A	501	4S1	C27-C26	2.95	1.43	1.36
2	D	501	4S1	C09-N25	3.00	1.35	1.31
2	B	501	4S1	C05-C06	3.05	1.44	1.39
2	C	501	4S1	C09-N25	3.13	1.36	1.31
2	A	501	4S1	C05-C06	3.13	1.45	1.39
2	A	501	4S1	C28-C29	3.14	1.43	1.36
2	B	501	4S1	C09-N25	3.31	1.36	1.31
2	C	501	4S1	C09-S22	3.38	1.78	1.73
2	D	501	4S1	C03-N08	3.47	1.40	1.34
2	D	501	4S1	C03-N04	3.51	1.39	1.34
2	C	501	4S1	C03-N08	3.71	1.40	1.34
2	A	501	4S1	C03-N08	3.79	1.40	1.34
2	D	501	4S1	C05-C06	3.83	1.46	1.39
2	C	501	4S1	C10-N07	4.19	1.50	1.45
2	B	501	4S1	C09-S22	4.67	1.80	1.73

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	4S1	C06-C05-N04	-2.87	120.01	124.44
2	D	501	4S1	O18-C17-C13	-2.64	105.45	111.38
2	C	501	4S1	N07-C01-N02	-2.53	111.78	118.44
2	B	501	4S1	C06-C05-N04	-2.50	120.58	124.44
2	B	501	4S1	O18-C17-C13	-2.47	105.83	111.38
2	C	501	4S1	N04-C03-N02	-2.14	124.38	126.66
2	B	501	4S1	C14-C13-C12	2.05	106.91	102.12
2	B	501	4S1	C13-C14-C10	2.10	105.68	103.45
2	B	501	4S1	C01-N07-C10	2.13	127.93	124.34
2	C	501	4S1	O16-C12-C11	2.15	118.82	111.86
2	D	501	4S1	C14-C10-N07	2.18	115.15	112.48
2	D	501	4S1	C01-N07-C10	2.25	128.13	124.34
2	C	501	4S1	C27-C26-C23	2.35	122.45	118.94
2	C	501	4S1	C23-C24-N25	2.49	113.88	108.12
2	A	501	4S1	C23-C24-N25	2.54	114.01	108.12
2	A	501	4S1	C01-N07-C10	2.58	128.70	124.34
2	B	501	4S1	C23-C24-N25	2.73	114.44	108.12
2	D	501	4S1	C13-C14-C10	2.77	106.39	103.45
2	D	501	4S1	C23-C24-N25	2.79	114.58	108.12
2	D	501	4S1	C09-N25-C24	3.26	110.22	103.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	4S1	C09-N25-C24	3.45	110.60	103.83
2	A	501	4S1	C13-C14-C10	3.47	107.14	103.45
2	A	501	4S1	C13-C12-C11	3.71	107.09	103.53
2	B	501	4S1	C13-C12-C11	3.74	107.12	103.53
2	B	501	4S1	C09-N25-C24	3.77	111.22	103.83
2	C	501	4S1	C19-N08-C03	3.82	130.13	123.62
2	B	501	4S1	C19-N08-C03	3.92	130.31	123.62
2	A	501	4S1	C09-N25-C24	4.07	111.82	103.83
2	D	501	4S1	C19-N08-C03	4.41	131.15	123.62
2	D	501	4S1	C13-C12-C11	4.44	107.79	103.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	4S1	2	0
2	B	501	4S1	2	0
2	C	501	4S1	3	0
2	D	501	4S1	3	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	280/301 (93%)	1.47	62 (22%) 1 1	39, 73, 148, 178	0
1	B	269/301 (89%)	1.33	58 (21%) 1 1	37, 69, 145, 162	0
1	C	282/301 (93%)	1.05	41 (14%) 3 3	33, 64, 126, 155	0
1	D	285/301 (94%)	0.82	32 (11%) 7 7	35, 63, 111, 153	0
All	All	1116/1204 (92%)	1.17	193 (17%) 2 2	33, 66, 141, 178	0

All (193) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	197	PHE	16.2
1	A	216	ALA	15.5
1	B	240	CYS	15.1
1	A	240	CYS	14.5
1	A	237	MET	10.8
1	B	187	VAL	10.7
1	A	223	THR	10.6
1	C	196	GLY	9.9
1	B	188	GLY	9.2
1	A	226	LEU	9.2
1	B	204	TYR	9.1
1	D	217	ALA	8.9
1	A	230	PHE	8.7
1	A	197	PHE	8.6
1	B	257	ASP	8.4
1	A	165	PHE	8.3
1	A	215	LEU	7.6
1	B	237	MET	7.6
1	A	229	GLN	7.4
1	B	216	ALA	7.2
1	C	195	GLY	6.8

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Mol	Chain	Res	Type	RSRZ
1	B	331	GLY	6.7
1	A	222	THR	6.6
1	C	226	LEU	6.5
1	A	236	VAL	6.4
1	A	204	TYR	6.4
1	C	191	LYS	6.3
1	C	215	LEU	6.2
1	A	227	LYS	6.2
1	B	185	ILE	6.2
1	A	188	GLY	6.2
1	B	215	LEU	6.2
1	A	258	LEU	6.0
1	A	195	GLY	5.9
1	C	187	VAL	5.5
1	C	230	PHE	5.4
1	A	334	ARG	5.4
1	A	185	ILE	5.4
1	B	186	SER	5.4
1	D	216	ALA	5.3
1	B	347	ARG	5.3
1	C	258	LEU	5.3
1	D	223	THR	5.3
1	A	256	ASP	5.2
1	B	333	ALA	5.2
1	D	187	VAL	5.2
1	D	230	PHE	5.2
1	D	256	ASP	5.0
1	A	239	LYS	4.9
1	A	225	GLU	4.8
1	A	254	ASP	4.7
1	C	223	THR	4.7
1	A	171	TYR	4.7
1	B	168	PHE	4.5
1	A	187	VAL	4.5
1	C	180	PHE	4.4
1	A	331	GLY	4.4
1	C	214	LYS	4.4
1	A	253	SER	4.3
1	B	198	GLY	4.3
1	D	186	SER	4.3
1	B	332	LEU	4.2
1	A	183	ARG	4.2

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Mol	Chain	Res	Type	RSRZ
1	B	330	PHE	4.2
1	B	236	VAL	4.2
1	D	226	LEU	4.1
1	B	256	ASP	4.1
1	A	407	GLU	4.1
1	C	183	ARG	4.1
1	B	191	LYS	4.1
1	C	185	ILE	4.1
1	B	334	ARG	4.0
1	B	184	PRO	4.0
1	A	228	GLN	4.0
1	A	251	PHE	3.9
1	C	227	LYS	3.9
1	B	233	GLU	3.8
1	A	238	ALA	3.8
1	B	253	SER	3.8
1	A	234	ILE	3.7
1	D	227	LYS	3.7
1	A	205	VAL	3.7
1	D	236	VAL	3.7
1	B	165	PHE	3.7
1	B	239	LYS	3.7
1	C	229	GLN	3.6
1	D	197	PHE	3.6
1	C	231	ASP	3.6
1	C	410	ILE	3.6
1	A	248	LEU	3.5
1	D	224	GLU	3.5
1	A	186	SER	3.5
1	A	196	GLY	3.5
1	D	349	VAL	3.4
1	B	173	LEU	3.4
1	B	234	ILE	3.4
1	C	236	VAL	3.4
1	C	192	MET	3.4
1	A	224	GLU	3.3
1	D	337	GLU	3.3
1	B	166	HIS	3.3
1	D	204	TYR	3.3
1	A	166	HIS	3.3
1	B	260	LEU	3.2
1	B	343	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	222	THR	3.2
1	D	195	GLY	3.2
1	C	171	TYR	3.2
1	B	251	PHE	3.2
1	C	343	VAL	3.2
1	A	191	LYS	3.2
1	B	189	GLY	3.2
1	C	332	LEU	3.2
1	C	186	SER	3.1
1	C	221	ILE	3.1
1	A	184	PRO	3.1
1	D	343	VAL	3.1
1	A	189	GLY	3.0
1	B	171	TYR	3.0
1	A	332	LEU	3.0
1	C	213	LYS	3.0
1	B	214	LYS	2.9
1	D	229	GLN	2.9
1	A	252	SER	2.9
1	B	183	ARG	2.9
1	D	347	ARG	2.9
1	A	333	ALA	2.9
1	A	330	PHE	2.8
1	A	257	ASP	2.8
1	D	239	LYS	2.8
1	C	204	TYR	2.8
1	C	190	ASN	2.7
1	A	399	ILE	2.7
1	B	407	GLU	2.7
1	C	167	SER	2.7
1	B	212	VAL	2.6
1	B	190	ASN	2.6
1	D	407	GLU	2.6
1	B	195	GLY	2.6
1	B	205	VAL	2.6
1	D	237	MET	2.6
1	D	254	ASP	2.6
1	A	206	ASN	2.6
1	B	348	ILE	2.6
1	B	199	VAL	2.5
1	A	347	ARG	2.5
1	A	207	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	364	ILE	2.5
1	C	290	LYS	2.5
1	B	194	GLU	2.5
1	B	254	ASP	2.5
1	A	260	LEU	2.4
1	A	214	LYS	2.4
1	C	239	LYS	2.4
1	B	252	SER	2.4
1	A	170	PHE	2.4
1	A	343	VAL	2.4
1	B	301	PHE	2.4
1	D	183	ARG	2.4
1	D	255	GLY	2.4
1	C	237	MET	2.4
1	A	246	VAL	2.4
1	D	215	LEU	2.4
1	A	348	ILE	2.3
1	A	190	ASN	2.3
1	D	185	ILE	2.3
1	B	241	GLN	2.3
1	C	165	PHE	2.3
1	C	211	ALA	2.3
1	A	247	GLU	2.2
1	B	341	GLN	2.2
1	B	360	LEU	2.2
1	A	443	LYS	2.2
1	D	443	LYS	2.2
1	B	235	LYS	2.2
1	D	460	SER	2.2
1	C	228	GLN	2.2
1	B	349	VAL	2.1
1	D	240	CYS	2.1
1	B	261	VAL	2.1
1	A	180	PHE	2.1
1	B	206	ASN	2.1
1	C	344	MET	2.1
1	B	443	LYS	2.1
1	C	452	GLN	2.1
1	C	177	THR	2.1
1	D	432	VAL	2.1
1	B	245	LEU	2.1
1	C	234	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	188	GLY	2.0
1	B	258	LEU	2.0
1	B	170	PHE	2.0
1	C	170	PHE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	SEP	D	346	10/11	0.63	0.24	-	90,95,103,103	0
1	SEP	C	346	10/11	0.68	0.18	-	88,94,102,102	0
1	SEP	B	346	10/11	0.68	0.23	-	108,112,117,118	0
1	SEP	A	346	10/11	0.78	0.22	-	107,110,117,117	0
1	TPO	A	342	11/12	0.63	0.26	-	114,116,120,120	0
1	TPO	B	342	11/12	0.88	0.29	-	111,113,116,117	0
1	TPO	B	345	11/12	0.91	0.13	-	104,104,108,109	0
1	TPO	C	342	11/12	0.84	0.17	-	99,100,106,106	0
1	TPO	C	345	11/12	0.88	0.16	-	86,87,89,90	0
1	TPO	A	345	11/12	0.90	0.14	-	102,103,108,109	0
1	TPO	D	345	11/12	0.92	0.13	-	86,89,91,91	0
1	TPO	D	342	11/12	0.84	0.17	-	98,100,107,107	0

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	4S1	A	501	29/29	0.88	0.21	0.19	75,77,82,84	0
2	4S1	C	501	29/29	0.89	0.22	0.18	63,71,79,80	0
2	4S1	B	501	29/29	0.90	0.18	-0.32	70,73,81,88	0
2	4S1	D	501	29/29	0.95	0.16	-0.32	54,60,72,75	0

6.5 Other polymers [i](#)

There are no such residues in this entry.