



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:12 PM UTC

PDB ID : 3KUK / pdb_00003kuk
Title : Trapping of an oxocarbenium ion intermediate in UP crystals
Authors : Paul, D.; O'Leary, S.; Rajashankar, K.; Bu, W.; Toms, A.; Settembre, E.;
Sanders, J.; Begley, T.P.; Ealick, S.E.
Deposited on : 2009-11-27
Resolution : 2.78 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

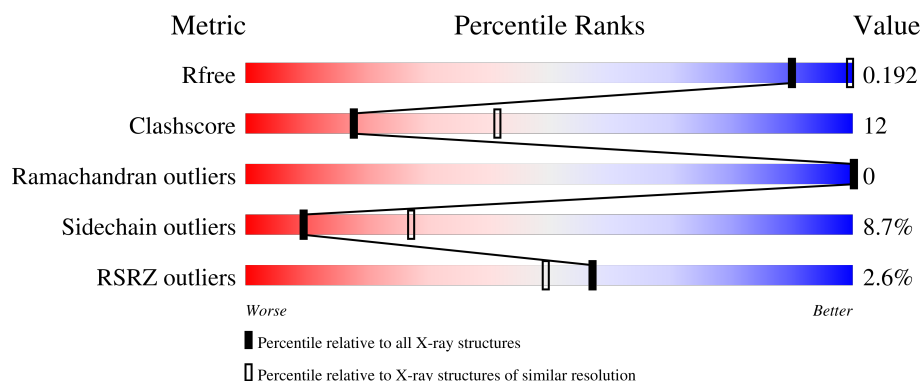
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

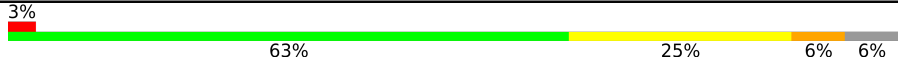
The reported resolution of this entry is 2.78 Å.

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	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	309	
1	B	309	

2 Entry composition [i](#)

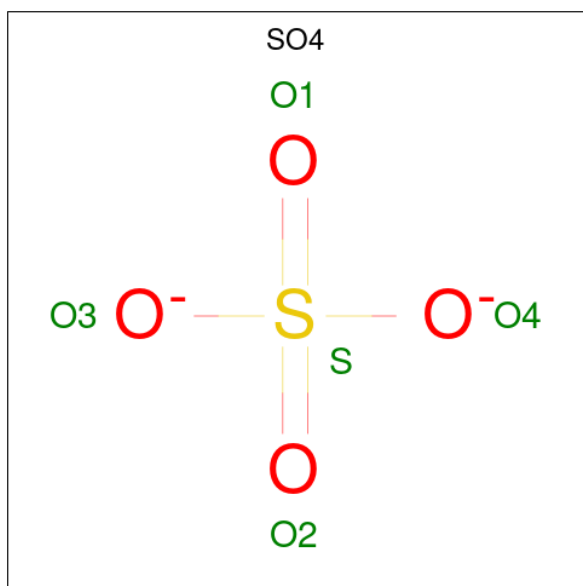
There are 4 unique types of molecules in this entry. The entry contains 4586 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Uridine phosphorylase.

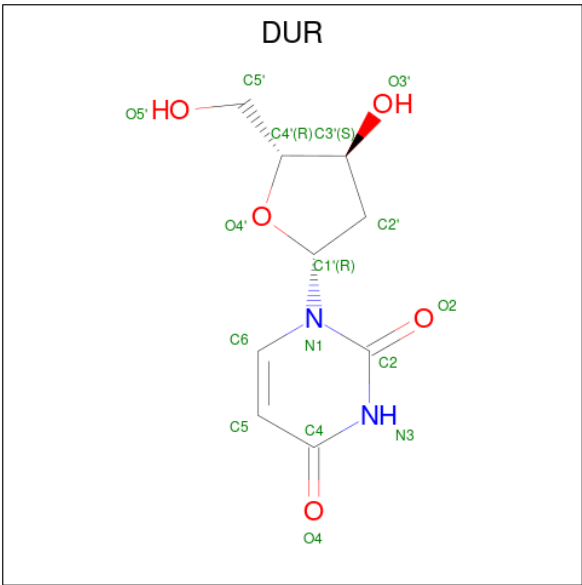
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	290	Total	C	N	O	S	0	0	0
			2226	1408	382	414	22			
1	B	293	Total	C	N	O	S	0	0	0
			2258	1427	391	418	22			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 2'-DEOXYURIDINE (CCD ID: DUR) (formula: C₉H₁₂N₂O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			16	9	2	5		
3	B	1	Total	C	N	O	0	0
			16	9	2	5		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	31	Total	O	0	0
			31	31		
4	B	29	Total	O	0	0
			29	29		

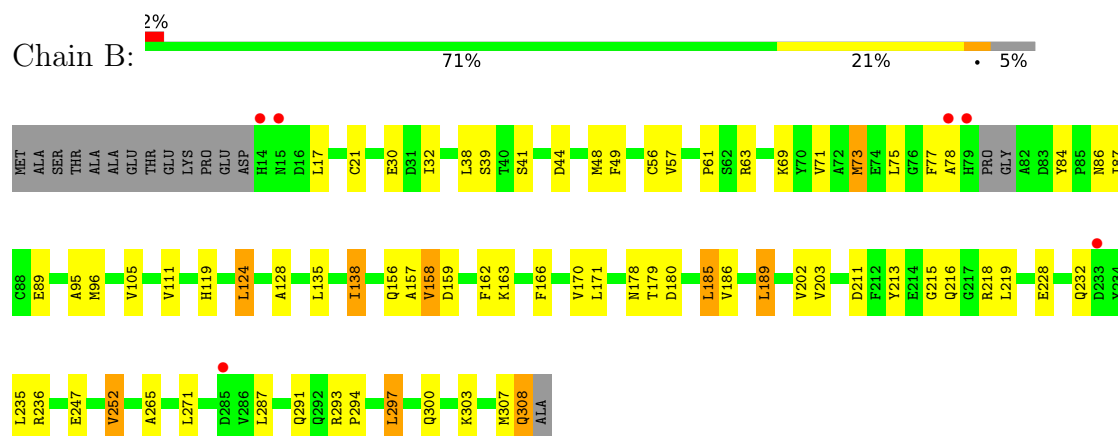
3 Residue-property plots [i](#)

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: Uridine phosphorylase



• Molecule 1: Uridine phosphorylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	82.64Å 82.64Å 259.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.78 78.73 – 2.78	Depositor EDS
% Data completeness (in resolution range)	83.5 ((Not available)-2.78) 88.2 (78.73-2.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.77Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.182 , 0.250 0.195 , 0.192	Depositor DCC
R_{free} test set	1052 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	34.3	Xtriage
Anisotropy	0.857	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 30.2	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	0.94	EDS
Total number of atoms	4586	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.95% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: DUR, SO4

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	0.45	0/2267	0.89	4/3063 (0.1%)
1	B	0.44	0/2301	0.82	0/3108
All	All	0.44	0/4568	0.86	4/6171 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	22	ASN	CA-C-N	6.56	126.49	119.28
1	A	22	ASN	C-N-CA	6.56	126.49	119.28
1	A	84	TYR	CA-C-N	6.27	126.49	119.90
1	A	84	TYR	C-N-CA	6.27	126.49	119.90

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	2226	0	2214	61	0
1	B	2258	0	2245	53	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	16	0	12	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	16	0	12	2	0
4	A	31	0	0	1	0
4	B	29	0	0	0	0
All	All	4586	0	4483	111	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 12.

All (111) 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:223:LEU:HD12	1:B:21:CYS:HB2	1.61	0.81
1:B:178:ASN:HD21	1:B:180:ASP:HB2	1.49	0.77
1:A:107:HIS:HB3	1:A:117:MET:HE2	1.69	0.75
3:A:311:DUR:O2	3:A:311:DUR:H2'1	1.87	0.74
1:B:56:CYS:HB3	1:B:138:ILE:HD11	1.70	0.73
1:B:232:GLN:O	1:B:236:ARG:HG3	1.89	0.73
1:A:279:GLN:HE22	1:B:32:ILE:H	1.34	0.73
1:A:138:ILE:O	1:A:138:ILE:HG12	1.90	0.70
1:A:185:LEU:HD13	1:A:265:ALA:HB2	1.75	0.67
1:A:173:LYS:HD3	4:A:324:HOH:O	1.93	0.67
1:A:166:PHE:CZ	1:A:252:VAL:HG13	2.30	0.67
1:A:31:ASP:H	1:A:40:THR:HB	1.59	0.66
1:B:297:LEU:O	1:B:297:LEU:HD12	1.95	0.66
3:B:311:DUR:H2'1	3:B:311:DUR:O2	1.94	0.65
1:A:124:LEU:C	1:A:124:LEU:HD23	2.21	0.65
1:A:222:ALA:H	1:B:119:HIS:HD2	1.45	0.65
1:A:189:LEU:HD12	1:A:297:LEU:HD21	1.79	0.64
1:B:211:ASP:HB3	1:B:215:GLY:HA3	1.80	0.64
1:B:166:PHE:CZ	1:B:252:VAL:HG13	2.35	0.61
1:B:185:LEU:HD22	1:B:189:LEU:HD22	1.83	0.61
1:A:159:ASP:HB2	1:A:160:PRO:CD	2.32	0.60
1:B:307:MET:C	1:B:308:GLN:HG3	2.25	0.60
1:B:111:VAL:HG22	1:B:252:VAL:HG22	1.84	0.59
1:B:185:LEU:HD13	1:B:265:ALA:HB2	1.84	0.59
1:B:300:GLN:NE2	1:B:300:GLN:HA	2.18	0.58
1:A:107:HIS:HD2	1:A:108:GLY:O	1.86	0.58
1:A:189:LEU:HD12	1:A:297:LEU:CD2	2.33	0.57
1:A:220:ASP:OD1	1:A:279:GLN:HG2	2.05	0.57
1:A:55:VAL:HG11	1:A:121:LEU:HD21	1.87	0.57
1:A:38:LEU:HD12	1:A:43:HIS:CE1	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:LYS:HE3	1:A:84:TYR:CD2	2.39	0.57
1:B:271:LEU:HD22	1:B:287:LEU:HD22	1.88	0.56
1:A:73:MET:HG2	1:A:78:ALA:HB2	1.87	0.56
1:A:185:LEU:HD22	1:A:189:LEU:HD22	1.89	0.55
1:B:158:VAL:HG13	1:B:162:PHE:HA	1.88	0.55
1:B:297:LEU:HD12	1:B:297:LEU:C	2.32	0.54
1:B:69:LYS:HG2	1:B:84:TYR:CE2	2.42	0.54
1:B:124:LEU:C	1:B:124:LEU:HD23	2.33	0.54
1:A:39:SER:O	1:A:43:HIS:HB2	2.08	0.54
1:B:69:LYS:HE2	1:B:84:TYR:CD2	2.43	0.54
1:B:158:VAL:CG1	1:B:162:PHE:HA	2.37	0.53
1:A:303:LYS:HG2	1:A:307:MET:HE3	1.90	0.53
1:B:73:MET:HA	1:B:78:ALA:HB2	1.91	0.53
1:B:86:ASN:O	1:B:89:GLU:HG2	2.09	0.52
1:A:148:PRO:HA	1:A:270:THR:HG22	1.92	0.52
1:B:300:GLN:HA	1:B:300:GLN:HE21	1.75	0.52
1:A:182:ASP:OD2	1:A:263:ARG:HG2	2.10	0.52
1:B:30:GLU:HG3	1:B:41:SER:HB3	1.91	0.51
1:B:293:ARG:HB2	1:B:294:PRO:HD3	1.92	0.51
1:A:73:MET:HG2	1:A:78:ALA:CB	2.40	0.51
1:A:196:LEU:HG	1:A:300:GLN:HG3	1.92	0.51
1:A:269:VAL:HG21	1:A:290:TYR:HB3	1.92	0.50
1:A:107:HIS:CD2	1:A:108:GLY:O	2.64	0.50
1:A:182:ASP:O	1:A:186:VAL:HG13	2.12	0.49
1:A:71:VAL:HG12	1:A:72:ALA:N	2.28	0.49
1:A:36:PHE:O	1:A:37:SER:HB3	2.13	0.48
1:B:44:ASP:O	1:B:48:MET:HG3	2.12	0.48
1:A:212:PHE:CE1	1:A:218:ARG:CZ	2.96	0.48
1:B:158:VAL:HG22	1:B:163:LYS:C	2.39	0.47
1:A:57:VAL:HA	1:A:105:VAL:O	2.14	0.47
1:B:73:MET:HG3	1:B:78:ALA:HB2	1.95	0.47
1:A:296:ARG:HH11	1:A:296:ARG:HG2	1.80	0.47
1:A:146:LEU:HD11	1:A:202:VAL:HG21	1.97	0.46
1:A:158:VAL:HG22	1:A:163:LYS:C	2.41	0.46
1:B:73:MET:HG3	1:B:78:ALA:CB	2.46	0.46
1:B:211:ASP:O	1:B:216:GLN:HG3	2.15	0.46
1:A:279:GLN:NE2	1:B:32:ILE:H	2.10	0.46
1:B:178:ASN:ND2	1:B:180:ASP:HB2	2.25	0.46
1:A:65:LYS:HG3	1:A:97:PHE:CE2	2.51	0.45
1:A:159:ASP:HB2	1:A:160:PRO:HD3	1.98	0.45
1:A:207:MET:HE2	1:A:246:ILE:HG12	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:300:GLN:HE21	1:B:300:GLN:CA	2.30	0.45
1:A:35:HIS:NE2	3:B:311:DUR:H5'2	2.32	0.45
1:A:22:ASN:HA	1:A:23:PRO:HD2	1.74	0.44
1:A:297:LEU:C	1:A:297:LEU:HD23	2.42	0.44
1:A:28:MET:HE2	1:A:30:GLU:O	2.18	0.44
1:B:170:VAL:HG12	1:B:171:LEU:HG	1.99	0.44
1:B:211:ASP:HB3	1:B:215:GLY:CA	2.46	0.44
1:A:135:LEU:O	1:A:264:ALA:HA	2.17	0.44
1:A:162:PHE:N	1:A:162:PHE:CD2	2.83	0.44
1:B:63:ARG:HD2	1:B:291:GLN:OE1	2.18	0.44
1:A:105:VAL:HG12	1:A:106:SER:O	2.18	0.44
1:A:124:LEU:HD23	1:A:125:LEU:N	2.32	0.44
1:B:49:PHE:HB2	1:B:128:ALA:CB	2.48	0.43
1:B:75:LEU:HB2	1:B:77:PHE:CD1	2.53	0.43
1:A:170:VAL:O	1:A:173:LYS:HD2	2.19	0.43
1:A:159:ASP:CB	1:A:160:PRO:CD	2.97	0.42
1:B:157:ALA:HB3	1:B:179:THR:HB	2.01	0.42
1:B:17:LEU:HD23	1:B:17:LEU:HA	1.86	0.42
1:B:57:VAL:HA	1:B:105:VAL:O	2.20	0.42
1:B:235:LEU:HD23	1:B:235:LEU:HA	1.79	0.42
1:B:303:LYS:HG2	1:B:307:MET:HE3	2.02	0.42
1:A:39:SER:HB3	1:A:42:THR:OG1	2.20	0.42
1:A:292:GLN:O	1:A:295:GLN:HB2	2.20	0.42
1:A:60:SER:O	1:A:64:MET:HG2	2.20	0.42
1:B:87:ILE:HD12	1:B:96:MET:HG2	2.02	0.42
1:A:84:TYR:CD2	1:A:84:TYR:N	2.88	0.41
1:B:232:GLN:OE1	1:B:236:ARG:NH1	2.53	0.41
1:B:159:ASP:OD2	1:B:159:ASP:C	2.63	0.41
1:A:96:MET:HE3	1:A:96:MET:HB2	1.85	0.41
1:A:92:ASP:OD1	1:B:61:PRO:HD2	2.20	0.41
1:A:293:ARG:HB2	1:A:294:PRO:HD3	2.02	0.41
1:B:156:GLN:HG3	1:B:180:ASP:OD1	2.21	0.41
1:B:219:LEU:HD11	1:B:228:GLU:HA	2.03	0.41
1:A:107:HIS:HB2	1:A:114:ILE:HA	2.01	0.41
1:B:213:TYR:O	1:B:218:ARG:HB2	2.21	0.41
1:A:304:LYS:O	1:A:308:GLN:HG3	2.21	0.40
1:B:178:ASN:HD21	1:B:180:ASP:CB	2.23	0.40
1:A:48:MET:HE2	1:A:49:PHE:CZ	2.56	0.40
1:A:211:ASP:O	1:A:216:GLN:HG3	2.22	0.40
1:B:61:PRO:HB3	1:B:95:ALA:CB	2.52	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	286/309 (93%)	273 (96%)	13 (4%)	0	100	100
1	B	289/309 (94%)	278 (96%)	11 (4%)	0	100	100
All	All	575/618 (93%)	551 (96%)	24 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	240/255 (94%)	215 (90%)	25 (10%)	7	20
1	B	244/255 (96%)	227 (93%)	17 (7%)	14	37
All	All	484/510 (95%)	442 (91%)	42 (9%)	9	27

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
1	A	38	LEU
1	A	40	THR
1	A	62	SER
1	A	71	VAL
1	A	124	LEU
1	A	135	LEU
1	A	138	ILE

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Mol	Chain	Res	Type
1	A	140	THR
1	A	147	GLU
1	A	158	VAL
1	A	173	LYS
1	A	176	VAL
1	A	185	LEU
1	A	186	VAL
1	A	189	LEU
1	A	203	VAL
1	A	223	LEU
1	A	229	LYS
1	A	247	GLU
1	A	252	VAL
1	A	270	THR
1	A	280	ILE
1	A	281	SER
1	A	304	LYS
1	B	38	LEU
1	B	39	SER
1	B	71	VAL
1	B	73	MET
1	B	124	LEU
1	B	135	LEU
1	B	138	ILE
1	B	158	VAL
1	B	185	LEU
1	B	186	VAL
1	B	189	LEU
1	B	202	VAL
1	B	203	VAL
1	B	247	GLU
1	B	252	VAL
1	B	297	LEU
1	B	308	GLN

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	24	HIS
1	A	107	HIS
1	A	168	GLN
1	A	187	GLN

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Mol	Chain	Res	Type
1	A	258	ASN
1	A	279	GLN
1	B	14	HIS
1	B	79	HIS
1	B	107	HIS
1	B	119	HIS
1	B	178	ASN
1	B	184	GLN
1	B	300	GLN

5.3.3 RNA [i](#)

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	SO4	A	310	-	4,4,4	0.25	0	6,6,6	0.34	0
2	SO4	B	310	-	4,4,4	0.22	0	6,6,6	0.19	0
3	DUR	B	311	-	17,17,17	2.92	5 (29%)	24,24,24	2.11	9 (37%)
3	DUR	A	311	-	17,17,17	2.84	4 (23%)	24,24,24	2.02	6 (25%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	DUR	A	311	-	-	3/6/18/18	0/2/2/2
3	DUR	B	311	-	-	2/6/18/18	0/2/2/2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	311	DUR	O4-C4	8.53	1.41	1.24
3	A	311	DUR	O4-C4	8.46	1.41	1.24
3	B	311	DUR	C2-N1	5.65	1.47	1.38
3	A	311	DUR	C2-N1	5.43	1.47	1.38
3	B	311	DUR	C6-C5	4.08	1.44	1.35
3	A	311	DUR	C6-C5	3.74	1.43	1.35
3	A	311	DUR	C3'-C4'	2.83	1.60	1.53
3	B	311	DUR	C3'-C4'	2.76	1.60	1.53
3	B	311	DUR	O2-C2	-2.23	1.19	1.23

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	311	DUR	C4-N3-C2	-4.78	120.67	126.61
3	A	311	DUR	C4-N3-C2	-4.47	121.07	126.61
3	B	311	DUR	N3-C2-N1	4.08	120.20	114.89
3	A	311	DUR	C5-C4-N3	3.53	119.75	114.80
3	B	311	DUR	C5-C4-N3	3.44	119.61	114.80
3	B	311	DUR	O4'-C4'-C5'	3.26	116.10	109.22
3	B	311	DUR	O4'-C1'-N1	3.15	113.46	107.86
3	A	311	DUR	N3-C2-N1	3.15	118.99	114.89
3	A	311	DUR	O4'-C1'-N1	3.05	113.28	107.86
3	A	311	DUR	C1'-N1-C2	2.82	123.18	117.66
3	A	311	DUR	O4-C4-C5	-2.70	120.50	125.16
3	B	311	DUR	C1'-N1-C2	2.19	121.94	117.66
3	B	311	DUR	C5'-C4'-C3'	2.17	120.23	114.82
3	B	311	DUR	C3'-C2'-C1'	2.10	107.74	102.60
3	B	311	DUR	O4-C4-C5	-2.10	121.55	125.16

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	311	DUR	C2'-C1'-N1-C2
3	A	311	DUR	C2'-C1'-N1-C6
3	B	311	DUR	C2'-C1'-N1-C2
3	B	311	DUR	C2'-C1'-N1-C6
3	A	311	DUR	O4'-C4'-C5'-O5'

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	311	DUR	2	0
3	A	311	DUR	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 [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	290/309 (93%)	-0.14	9 (3%)	51 44	24, 36, 58, 74	0
1	B	293/309 (94%)	-0.25	6 (2%)	65 58	24, 36, 57, 99	0
All	All	583/618 (94%)	-0.19	15 (2%)	57 49	24, 36, 58, 99	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	233	ASP	5.3
1	B	14	HIS	4.5
1	B	285	ASP	4.1
1	B	233	ASP	4.0
1	B	15	ASN	3.3
1	B	79	HIS	3.1
1	A	307	MET	3.0
1	A	82	ALA	2.9
1	A	77	PHE	2.7
1	A	17	LEU	2.3
1	A	78	ALA	2.3
1	B	78	ALA	2.3
1	A	16	ASP	2.2
1	A	90	GLY	2.1
1	A	21	CYS	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
3	DUR	B	311	16/16	0.88	0.14	34,40,50,52	0
3	DUR	A	311	16/16	0.94	0.10	26,38,44,46	0
2	SO4	A	310	5/5	0.98	0.06	26,30,36,39	0
2	SO4	B	310	5/5	0.98	0.07	32,34,36,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.