



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 10, 2020 – 12:14 PM JST

PDB ID : 6J6T
Title : Crystal Structure of HDA15 HD domain
Authors : Cheng, Y.S.; Hsu, J.C.; Hung, H.C.; Liu, T.C.
Deposited on : 2019-01-15
Resolution : 2.36 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.14.6
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.14.6

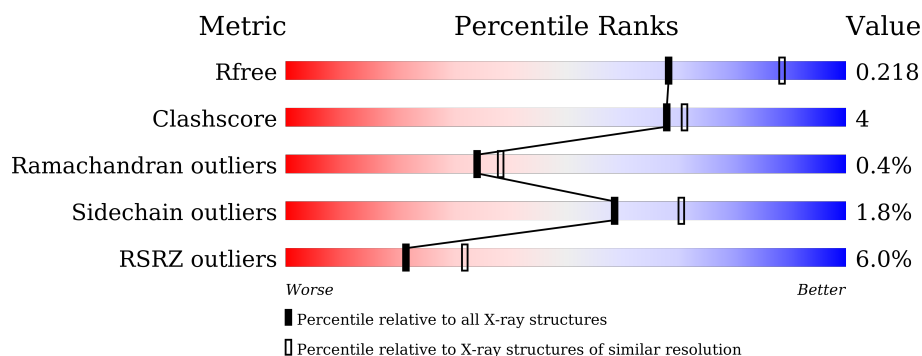
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.36 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}	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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 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	364	
1	B	364	
1	C	364	
1	D	364	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 11022 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 Histone deacetylase 15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	353	Total	C	N	O	S	0	0	0
			2647	1676	460	497	14			
1	B	356	Total	C	N	O	S	0	0	0
			2668	1685	463	506	14			
1	C	344	Total	C	N	O	S	0	0	0
			2586	1637	451	484	14			
1	D	351	Total	C	N	O	S	0	0	0
			2634	1664	458	498	14			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	K	0	0
			1	1		
3	A	1	Total	K	0	0
			1	1		
3	D	1	Total	K	0	0
			1	1		
3	C	1	Total	K	0	0
			1	1		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	139	Total	O	0	0
			139	139		
5	B	126	Total	O	0	0
			126	126		
5	C	51	Total	O	0	0
			51	51		

Continued on next page...

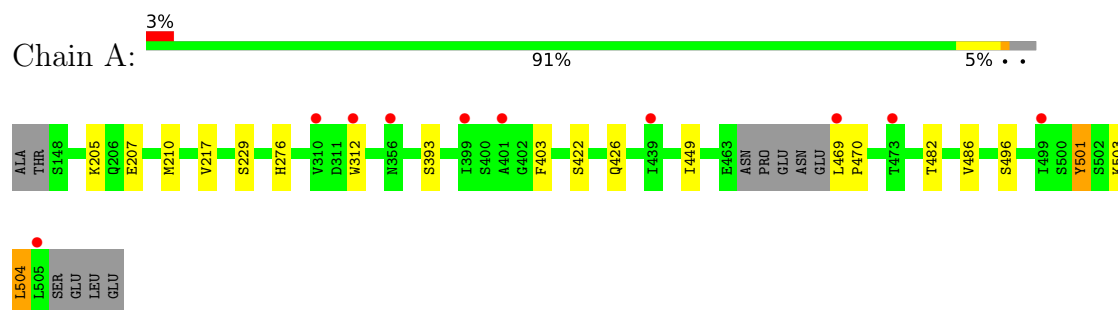
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	123	Total 123	O 123	0	0

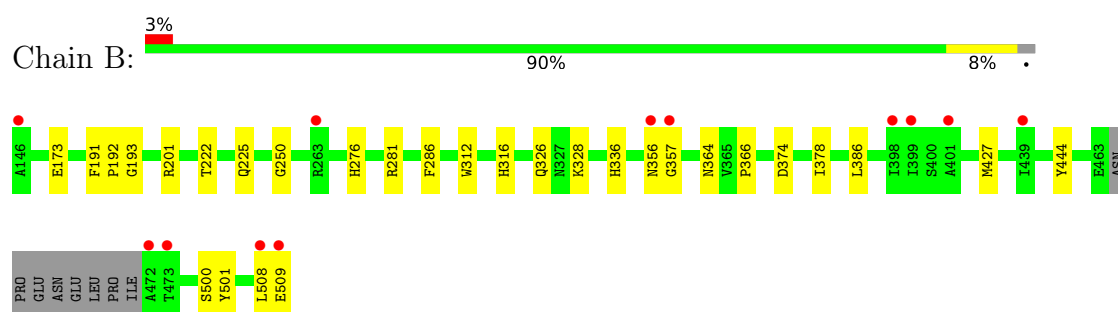
3 Residue-property plots

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.

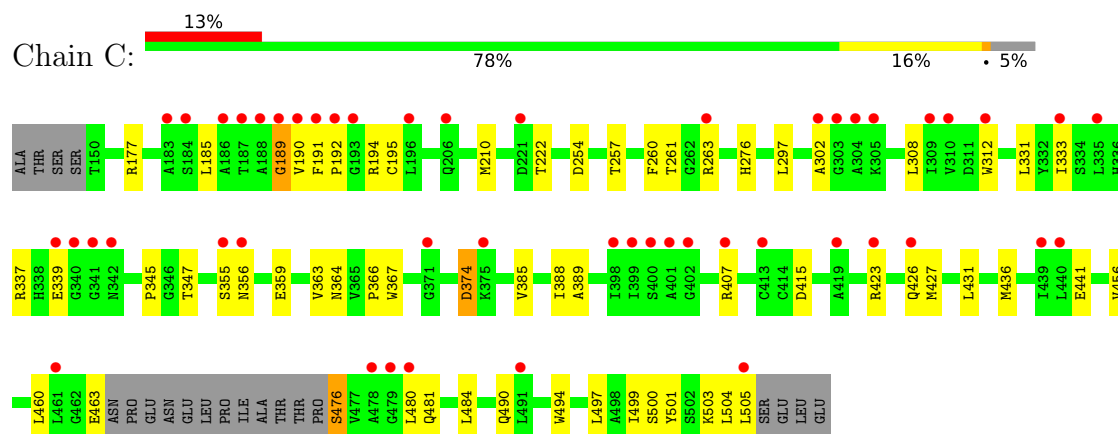
• Molecule 1: Histone deacetylase 15



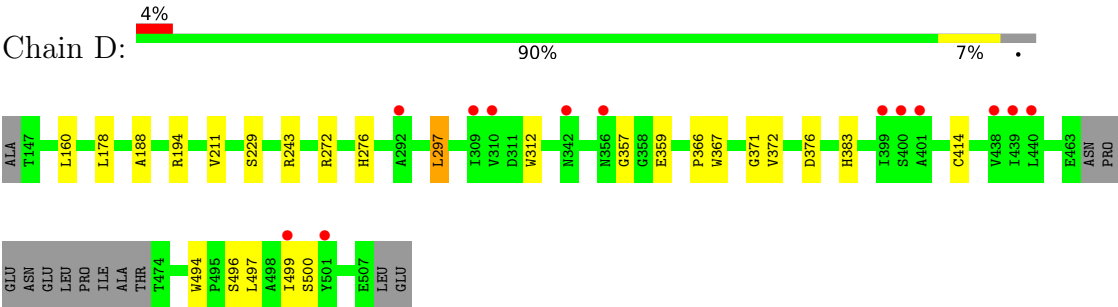
• Molecule 1: Histone deacetylase 15



• Molecule 1: Histone deacetylase 15



• Molecule 1: Histone deacetylase 15



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.81Å 78.83Å 176.79Å 90.00° 94.22° 90.00°	Depositor
Resolution (Å)	29.52 – 2.36 29.55 – 2.36	Depositor EDS
% Data completeness (in resolution range)	96.4 (29.52-2.36) 96.2 (29.55-2.36)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.27 (at 2.36Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.167 , 0.218 0.167 , 0.218	Depositor DCC
R_{free} test set	2001 reflections (2.92%)	wwPDB-VP
Wilson B-factor (Å ²)	22.6	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 49.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11022	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 3.41% 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: ZN, K, 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	0.44	0/2706	0.60	0/3680
1	B	0.45	0/2726	0.58	0/3706
1	C	0.39	0/2643	0.58	0/3591
1	D	0.41	0/2692	0.58	1/3659 (0.0%)
All	All	0.42	0/10767	0.59	1/14636 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	297	LEU	CA-CB-CG	-5.28	103.15	115.30

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	2647	0	2597	9	0
1	B	2668	0	2608	17	0
1	C	2586	0	2532	44	0
1	D	2634	0	2573	12	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	10	0	0	0	0
4	B	15	0	0	1	0
4	C	5	0	0	0	0
4	D	10	0	0	0	0
5	A	139	0	0	0	0
5	B	126	0	0	2	0
5	C	51	0	0	1	0
5	D	123	0	0	0	0
All	All	11022	0	10310	80	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 4.

All (80) 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:386:LEU:HD21	1:B:427:MET:HB3	1.78	0.66
1:C:389:ALA:HB1	1:C:431:LEU:HD21	1.81	0.62
1:D:178:LEU:HD13	1:D:272:ARG:HG3	1.83	0.59
1:C:388:ILE:HG12	1:C:497:LEU:HD11	1.85	0.58
1:C:190:VAL:HB	1:C:194:ARG:H	1.69	0.57
1:C:190:VAL:HG21	1:C:195:CYS:CB	2.35	0.56
1:A:501:TYR:C	1:A:503:LYS:H	2.08	0.56
1:D:494:TRP:HB2	1:D:497:LEU:HD22	1.88	0.56
1:C:463:GLU:N	1:C:463:GLU:OE1	2.40	0.55
1:A:422:SER:O	1:A:426:GLN:HG3	2.06	0.55
1:C:190:VAL:HG21	1:C:195:CYS:HB2	1.88	0.55
1:A:205:LYS:HG3	1:A:217:VAL:HG11	1.88	0.55
1:B:508:LEU:HD12	1:B:509:GLU:HB2	1.89	0.54
1:B:281:ARG:HG2	5:B:729:HOH:O	2.07	0.54
1:B:222:THR:HG21	1:C:222:THR:HB	1.90	0.54
1:B:225:GLN:NE2	5:B:703:HOH:O	2.42	0.53
1:A:207:GLU:O	1:A:210:MET:HG2	2.09	0.53
1:B:326:GLN:HG2	1:B:356:ASN:HB2	1.91	0.53
1:C:190:VAL:HG22	1:C:191:PHE:HB2	1.89	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:SER:HB3	1:D:229:SER:HB3	1.91	0.52
1:C:490:GLN:HG3	1:C:494:TRP:HD1	1.72	0.52
1:C:333:ILE:HG12	1:C:363:VAL:HB	1.92	0.51
1:C:185:LEU:HB3	1:C:191:PHE:H	1.76	0.51
1:C:190:VAL:HG21	1:C:195:CYS:SG	2.51	0.51
1:C:490:GLN:HG3	1:C:494:TRP:CD1	2.45	0.51
1:B:374:ASP:O	1:B:378:ILE:HG12	2.11	0.50
1:D:383:HIS:ND1	1:D:500:SER:OG	2.43	0.49
1:B:201:ARG:O	1:B:250:GLY:HA3	2.13	0.48
1:B:508:LEU:CD1	1:B:509:GLU:HB2	2.43	0.48
1:C:504:LEU:HA	1:C:505:LEU:C	2.34	0.47
1:C:190:VAL:HG22	1:C:191:PHE:CB	2.45	0.47
1:B:328:LYS:HG2	4:B:603:SO4:O2	2.14	0.47
1:C:210:MET:HB3	1:C:297:LEU:HD23	1.95	0.47
1:D:211:VAL:HG22	1:D:297:LEU:CD1	2.44	0.47
1:C:374:ASP:N	1:C:374:ASP:OD1	2.48	0.47
1:A:403:PHE:HB2	1:A:449:ILE:HG22	1.96	0.46
1:C:254:ASP:HA	1:C:257:THR:HG22	1.96	0.46
1:C:345:PRO:HG2	1:C:347:THR:OG1	2.16	0.46
1:C:481:GLN:HA	1:C:484:LEU:HG	1.98	0.45
1:C:190:VAL:HG22	1:C:191:PHE:CA	2.47	0.45
1:D:496:SER:O	1:D:499:ILE:HD12	2.17	0.45
1:C:456:VAL:O	1:C:460:LEU:HD12	2.18	0.44
1:C:185:LEU:O	1:C:191:PHE:N	2.51	0.44
1:B:378:ILE:HD12	1:B:427:MET:HE1	2.00	0.44
1:D:371:GLY:CA	1:D:414:CYS:HA	2.47	0.44
1:B:316:HIS:HB2	1:B:336:HIS:CD2	2.53	0.44
1:C:476:SER:N	5:C:704:HOH:O	2.49	0.44
1:C:480:LEU:HD23	1:C:484:LEU:HD23	2.00	0.44
1:D:188:ALA:O	1:D:194:ARG:NH2	2.50	0.43
1:D:160:LEU:HD12	1:D:243:ARG:NH2	2.33	0.43
1:A:469:LEU:HA	1:A:470:PRO:HD3	1.89	0.43
1:C:423:ARG:HH11	1:C:426:GLN:HB3	1.84	0.43
1:C:260:PHE:HE2	1:C:302:ALA:O	2.02	0.43
1:D:367:TRP:CG	1:D:372:VAL:HG21	2.54	0.43
1:C:423:ARG:NH1	1:C:426:GLN:HB3	2.34	0.42
1:C:190:VAL:HB	1:C:194:ARG:N	2.35	0.42
1:C:337:ARG:HA	1:C:367:TRP:HB2	2.01	0.42
1:C:261:THR:HG23	1:C:263:ARG:HG3	2.01	0.42
1:C:499:ILE:O	1:C:503:LYS:HG2	2.20	0.41
1:C:189:GLY:HA2	1:C:190:VAL:C	2.39	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:308:LEU:HD13	1:C:331:LEU:HD23	2.01	0.41
1:D:372:VAL:HG12	1:D:376:ASP:HB2	2.02	0.41
1:C:191:PHE:CD1	1:C:192:PRO:HA	2.55	0.41
1:C:356:ASN:O	1:C:359:GLU:HG3	2.20	0.41
1:C:333:ILE:HG21	1:C:385:VAL:HG22	2.01	0.41
1:C:388:ILE:CG1	1:C:497:LEU:HD11	2.50	0.41
1:B:364:ASN:O	1:B:366:PRO:HD3	2.21	0.41
1:C:190:VAL:CG2	1:C:191:PHE:HA	2.51	0.41
1:C:423:ARG:O	1:C:427:MET:HG3	2.21	0.41
1:C:436:MET:HB2	1:C:436:MET:HE2	1.89	0.41
1:B:328:LYS:HD2	1:B:357:GLY:O	2.20	0.41
1:A:482:THR:O	1:A:486:VAL:HG23	2.21	0.41
1:A:501:TYR:O	1:A:504:LEU:HD12	2.20	0.41
1:B:286:PHE:HD2	1:B:444:TYR:HH	1.67	0.40
1:C:190:VAL:CG2	1:C:195:CYS:H	2.35	0.40
1:C:177:ARG:NH2	1:C:441:GLU:O	2.51	0.40
1:D:357:GLY:N	1:D:359:GLU:OE2	2.51	0.40
1:B:191:PHE:HA	1:B:192:PRO:HA	1.95	0.40
1:B:192:PRO:HA	1:B:193:GLY:HA2	1.89	0.40
1:C:364:ASN:O	1:C:366:PRO:HD3	2.21	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	349/364 (96%)	334 (96%)	13 (4%)	2 (1%)	25	27
1	B	352/364 (97%)	338 (96%)	14 (4%)	0	100	100
1	C	340/364 (93%)	319 (94%)	19 (6%)	2 (1%)	25	27
1	D	347/364 (95%)	337 (97%)	9 (3%)	1 (0%)	41	47
All	All	1388/1456 (95%)	1328 (96%)	55 (4%)	5 (0%)	34	38

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	504	LEU
1	A	496	SER
1	C	189	GLY
1	C	355	SER
1	D	366	PRO

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	276/286 (96%)	272 (99%)	4 (1%)	67	78
1	B	278/286 (97%)	273 (98%)	5 (2%)	59	70
1	C	268/286 (94%)	259 (97%)	9 (3%)	37	46
1	D	275/286 (96%)	273 (99%)	2 (1%)	84	91
All	All	1097/1144 (96%)	1077 (98%)	20 (2%)	59	70

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

Mol	Chain	Res	Type
1	A	276	HIS
1	A	312	TRP
1	A	393	SER
1	A	501	TYR
1	B	173	GLU
1	B	276	HIS
1	B	312	TRP
1	B	500	SER
1	B	501	TYR
1	C	276	HIS
1	C	312	TRP
1	C	339	GLU
1	C	374	ASP
1	C	407	ARG
1	C	415	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	476	SER
1	C	500	SER
1	C	501	TYR
1	D	276	HIS
1	D	312	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	326	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 8 are monoatomic - leaving 8 for Mogul analysis.

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	$\# Z > 2$
4	SO4	B	605	-	4,4,4	0.18	0	6,6,6	0.27	0
4	SO4	B	604	-	4,4,4	0.14	0	6,6,6	0.25	0
4	SO4	B	603	-	4,4,4	0.18	0	6,6,6	0.14	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	A	603	-	4,4,4	0.12	0	6,6,6	0.36	0
4	SO4	D	603	-	4,4,4	0.23	0	6,6,6	0.18	0
4	SO4	D	604	-	4,4,4	0.31	0	6,6,6	0.38	0
4	SO4	C	603	-	4,4,4	0.12	0	6,6,6	0.12	0
4	SO4	A	604	-	4,4,4	0.14	0	6,6,6	0.22	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	603	SO4	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	353/364 (96%)	-0.18	10 (2%) 53 64	11, 20, 40, 65	0
1	B	356/364 (97%)	-0.16	12 (3%) 45 57	10, 22, 39, 74	0
1	C	344/364 (94%)	0.67	49 (14%) 2 4	15, 43, 64, 69	0
1	D	351/364 (96%)	-0.00	13 (3%) 41 54	13, 24, 40, 69	0
All	All	1404/1456 (96%)	0.08	84 (5%) 21 32	10, 24, 56, 74	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	187	THR	5.2
1	C	190	VAL	4.9
1	C	505	LEU	4.6
1	B	508	LEU	4.4
1	C	188	ALA	4.3
1	A	499	ILE	4.3
1	C	196	LEU	4.3
1	C	356	ASN	4.2
1	C	310	VAL	4.2
1	C	480	LEU	4.1
1	C	371	GLY	4.1
1	C	189	GLY	4.0
1	C	183	ALA	4.0
1	C	340	GLY	3.9
1	A	469	LEU	3.9
1	B	472	ALA	3.8
1	C	206	GLN	3.7
1	C	263	ARG	3.7
1	C	479	GLY	3.7
1	C	439	ILE	3.6
1	C	400	SER	3.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	401	ALA	3.5
1	C	192	PRO	3.5
1	C	341	GLY	3.5
1	C	303	GLY	3.5
1	C	186	ALA	3.5
1	D	356	ASN	3.4
1	C	342	ASN	3.4
1	C	423	ARG	3.4
1	B	146	ALA	3.4
1	C	309	ILE	3.3
1	D	439	ILE	3.3
1	D	401	ALA	3.2
1	C	407	ARG	3.1
1	C	399	ILE	3.1
1	D	499	ILE	3.1
1	C	312	TRP	3.0
1	B	399	ILE	3.0
1	B	509	GLU	3.0
1	A	473	THR	2.9
1	B	473	THR	2.8
1	C	355	SER	2.8
1	C	461	LEU	2.8
1	A	401	ALA	2.7
1	C	333	ILE	2.7
1	A	310	VAL	2.7
1	C	491	LEU	2.7
1	C	398	ILE	2.6
1	D	440	LEU	2.6
1	C	413	CYS	2.6
1	D	342	ASN	2.6
1	A	439	ILE	2.6
1	C	375	LYS	2.5
1	A	356	ASN	2.5
1	D	310	VAL	2.5
1	C	191	PHE	2.4
1	C	419	ALA	2.4
1	B	398	ILE	2.4
1	B	263	ARG	2.4
1	C	304	ALA	2.4
1	D	292	ALA	2.4
1	D	438	VAL	2.4
1	D	501	TYR	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	401	ALA	2.3
1	C	184	SER	2.3
1	C	426	GLN	2.3
1	D	309	ILE	2.3
1	C	305	LYS	2.2
1	B	357	GLY	2.2
1	C	478	ALA	2.2
1	A	505	LEU	2.2
1	C	221	ASP	2.2
1	B	356	ASN	2.2
1	C	339	GLU	2.2
1	C	402	GLY	2.2
1	D	400	SER	2.1
1	B	439	ILE	2.1
1	A	399	ILE	2.1
1	C	335	LEU	2.1
1	C	440	LEU	2.1
1	D	399	ILE	2.1
1	A	312	TRP	2.1
1	C	193	GLY	2.1
1	C	302	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 monosaccharides 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(\AA^2)	Q<0.9
4	SO4	D	604	5/5	0.87	0.35	46,49,56,59	0
4	SO4	A	604	5/5	0.93	0.35	43,53,65,67	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	SO4	B	603	5/5	0.95	0.29	47,51,54,56	0
4	SO4	B	604	5/5	0.95	0.30	43,48,55,63	0
4	SO4	A	603	5/5	0.96	0.15	38,39,46,56	0
2	ZN	A	601	1/1	0.97	0.07	36,36,36,36	0
2	ZN	C	601	1/1	0.97	0.07	45,45,45,45	0
4	SO4	C	603	5/5	0.97	0.15	36,48,53,64	0
4	SO4	D	603	5/5	0.97	0.15	37,38,53,58	0
2	ZN	D	601	1/1	0.98	0.09	36,36,36,36	0
3	K	C	602	1/1	0.98	0.11	33,33,33,33	0
2	ZN	B	601	1/1	0.98	0.06	36,36,36,36	0
3	K	D	602	1/1	0.99	0.13	20,20,20,20	0
4	SO4	B	605	5/5	0.99	0.10	28,29,30,39	0
3	K	B	602	1/1	1.00	0.12	17,17,17,17	0
3	K	A	602	1/1	1.00	0.10	17,17,17,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.