



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 06:28 AM UTC

PDB ID : 1VZ5 / pdb_00001vz5
Title : Succinate Complex of AtsK
Authors : Mueller, I.; Stueckl, A.C.; Uson, I.; Kertesz, M.
Deposited on : 2004-05-14
Resolution : 2.15 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

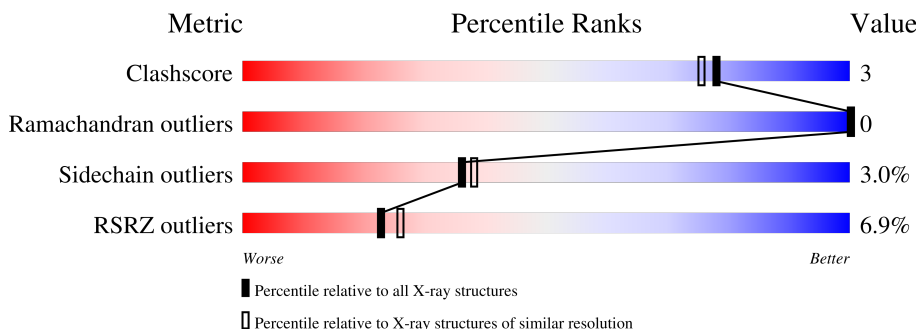
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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))
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	301	4% 76% 8% 16%
1	B	301	3% 73% 11% 16%
1	C	301	6% 71% 10% 19%
1	D	301	10% 70% 9% 19%

2 Entry composition [i](#)

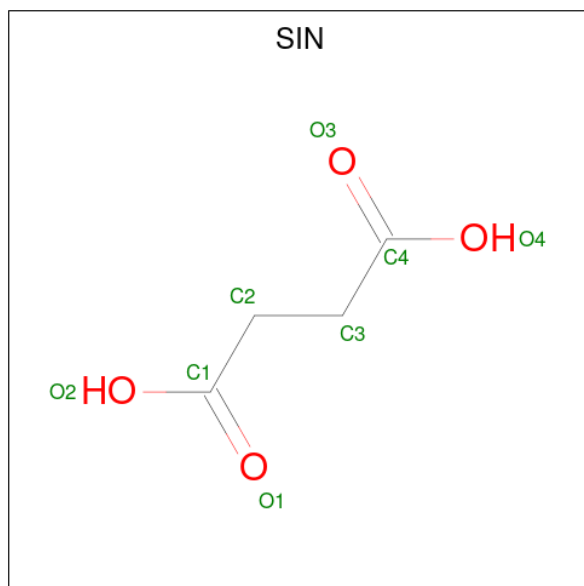
There are 3 unique types of molecules in this entry. The entry contains 8147 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PUTATIVE ALKYL SULFATASE ATSK.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	254	Total	C	N	O	0	0	0
			1959	1239	362	358			
1	B	254	Total	C	N	O	0	0	0
			1949	1233	356	360			
1	C	244	Total	C	N	O	0	0	0
			1841	1168	333	340			
1	D	244	Total	C	N	O	0	0	0
			1842	1170	334	338			

- Molecule 2 is SUCCINIC ACID (CCD ID: SIN) (formula: C₄H₆O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			8	4	4		
2	B	1	Total	C	O	0	0
			8	4	4		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			8	4	4		
2	D	1	Total	C	O	0	0
			8	4	4		

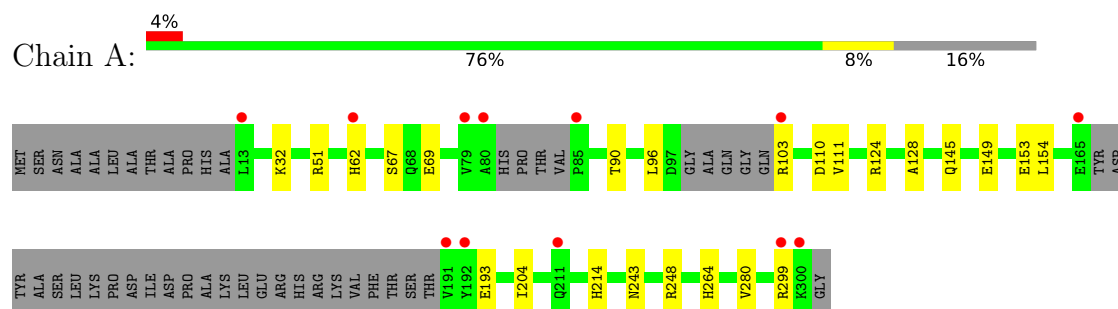
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	148	Total	O	0	0
			148	148		
3	B	139	Total	O	0	0
			139	139		
3	C	129	Total	O	0	0
			129	129		
3	D	108	Total	O	0	0
			108	108		

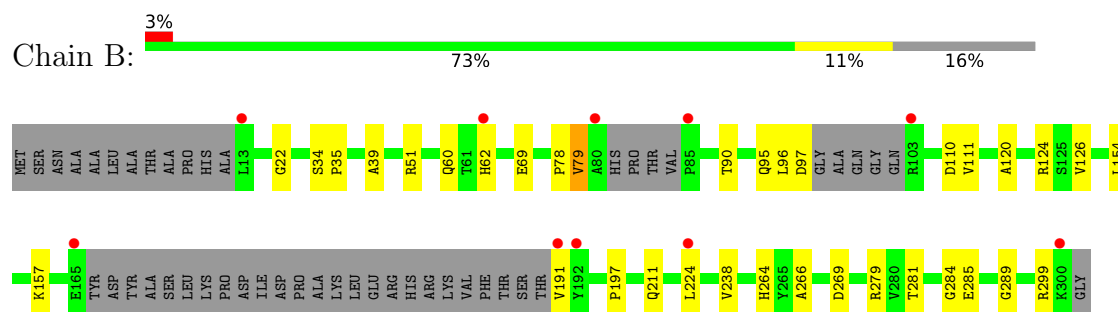
3 Residue-property plots

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.

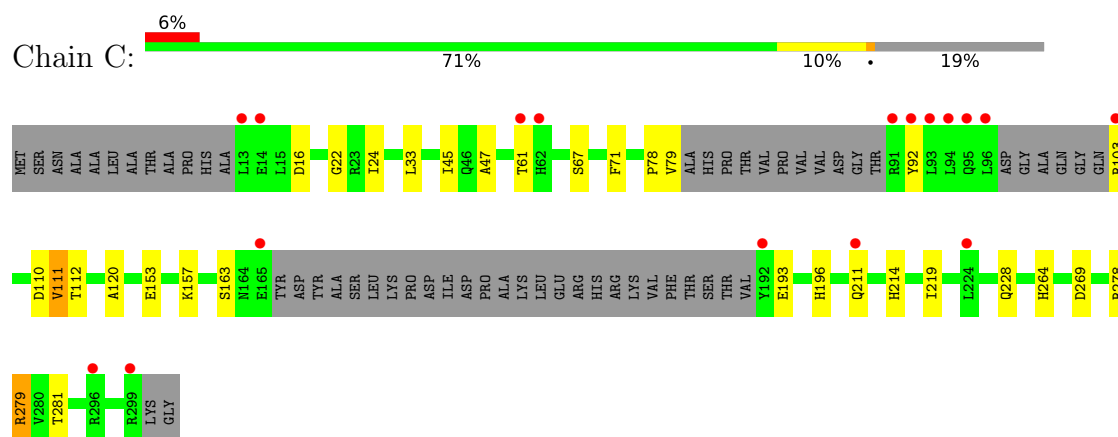
• Molecule 1: PUTATIVE ALKYL SULFATASE ATSK



• Molecule 1: PUTATIVE ALKYL SULFATASE ATSK

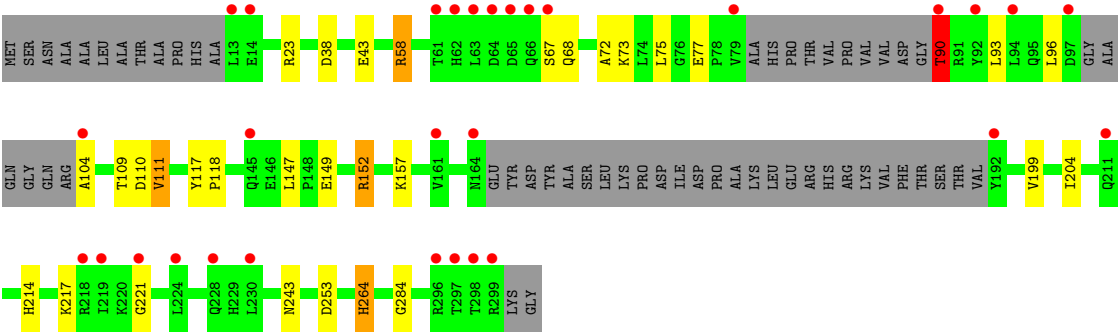


• Molecule 1: PUTATIVE ALKYL SULFATASE ATSK



• Molecule 1: PUTATIVE ALKYL SULFATASE ATSK





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.51Å 145.81Å 158.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.80 – 2.15 48.80 – 2.15	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.80-2.15) 100.0 (48.80-2.15)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.25 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
R, R_{free}	0.194 , 0.225 (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	25.6	Xtriage
Anisotropy	0.319	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 37.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8147	wwPDB-VP
Average B, all atoms (Å ²)	27.0	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 20.68 % 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 8.2399e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	1.45	6/2002 (0.3%)	1.19	1/2729 (0.0%)
1	B	1.46	11/1992 (0.6%)	1.22	1/2720 (0.0%)
1	C	1.48	9/1883 (0.5%)	1.21	1/2575 (0.0%)
1	D	1.58	10/1884 (0.5%)	1.30	9/2577 (0.3%)
All	All	1.49	36/7761 (0.5%)	1.23	12/10601 (0.1%)

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	149	GLU	CA-CB	10.41	1.66	1.53
1	C	278	ARG	N-CA	7.89	1.55	1.46
1	D	111	VAL	CA-CB	6.94	1.63	1.54
1	B	284	GLY	N-CA	6.91	1.50	1.44
1	A	204	ILE	CA-CB	6.82	1.62	1.55
1	C	111	VAL	CA-CB	6.56	1.63	1.54
1	A	67	SER	CA-C	6.54	1.61	1.52
1	C	24	ILE	CA-CB	6.54	1.63	1.54
1	D	284	GLY	N-CA	6.48	1.50	1.44
1	B	238	VAL	C-O	-6.29	1.17	1.24
1	B	120	ALA	N-CA	6.04	1.53	1.46
1	D	72	ALA	CA-CB	-5.93	1.43	1.53
1	D	152	ARG	N-CA	5.70	1.53	1.46
1	A	280	VAL	CA-CB	5.70	1.61	1.54
1	B	51	ARG	N-CA	-5.64	1.39	1.46
1	A	243	ASN	N-CA	5.59	1.53	1.46
1	D	67	SER	N-CA	-5.48	1.39	1.46
1	C	45	ILE	CA-C	5.46	1.59	1.52
1	A	128	ALA	CA-C	5.37	1.58	1.52
1	D	204	ILE	CA-CB	5.36	1.60	1.55
1	C	16	ASP	CA-C	5.30	1.59	1.52
1	D	199	VAL	CA-CB	-5.29	1.47	1.54

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	77	GLU	CA-CB	5.28	1.61	1.52
1	B	39	ALA	CA-CB	-5.24	1.44	1.53
1	D	264	HIS	C-O	5.22	1.30	1.23
1	B	126	VAL	C-O	5.21	1.29	1.23
1	D	75	LEU	N-CA	5.18	1.52	1.46
1	C	47	ALA	CA-CB	5.18	1.61	1.53
1	B	197	PRO	C-O	5.17	1.29	1.23
1	B	266	ALA	CA-CB	5.16	1.60	1.53
1	B	79	VAL	CA-CB	5.13	1.60	1.54
1	C	120	ALA	CA-C	-5.08	1.46	1.52
1	C	92	TYR	CA-C	5.06	1.59	1.52
1	C	211	GLN	CG-CD	5.05	1.64	1.52
1	B	289	GLY	C-O	5.03	1.29	1.23
1	B	224	LEU	N-CA	-5.00	1.40	1.46

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	58	ARG	NE-CZ-NH2	-7.07	112.84	119.20
1	D	73	LYS	N-CA-C	-6.81	104.84	113.01
1	D	204	ILE	CB-CA-C	-6.55	105.21	111.44
1	A	248	ARG	NE-CZ-NH1	-6.20	115.30	121.50
1	D	217	LYS	N-CA-C	-5.89	105.76	113.12
1	D	68	GLN	N-CA-C	-5.88	104.79	111.14
1	D	43	GLU	N-CA-C	-5.73	105.12	111.36
1	D	221	GLY	N-CA-C	-5.57	107.31	115.72
1	B	211	GLN	CB-CA-C	-5.42	102.11	111.23
1	C	67	SER	N-CA-CB	-5.38	102.20	110.16
1	D	90	THR	N-CA-C	5.17	125.47	111.00
1	D	109	THR	N-CA-C	-5.03	102.07	110.17

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1959	0	1916	9	0
1	B	1949	0	1887	14	0
1	C	1841	0	1746	11	0
1	D	1842	0	1748	13	0
2	A	8	0	4	0	0
2	B	8	0	4	0	0
2	C	8	0	4	0	0
2	D	8	0	4	0	0
3	A	148	0	0	2	0
3	B	139	0	0	0	0
3	C	129	0	0	0	0
3	D	108	0	0	2	0
All	All	8147	0	7313	42	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 (42) 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:23:ARG:NH2	1:D:243:ASN:OD1	1.86	1.07
1:C:153:GLU:OE1	1:D:157:LYS:NZ	2.16	0.73
1:D:58:ARG:HD3	1:D:253:ASP:OD1	1.93	0.69
1:A:193:GLU:HB2	1:A:299:ARG:HB3	1.78	0.66
1:A:110:ASP:OD1	1:A:264:HIS:HE1	1.80	0.64
1:D:38:ASP:C	1:D:38:ASP:OD1	2.41	0.62
1:D:90:THR:HG21	1:D:93:LEU:HB3	1.80	0.62
1:A:145:GLN:HG2	3:A:2090:HOH:O	2.01	0.59
1:A:103:ARG:NH2	3:A:2048:HOH:O	2.36	0.58
1:A:32:LYS:HG3	1:A:62:HIS:CE1	2.40	0.56
1:C:110:ASP:OD1	1:C:264:HIS:HE1	1.88	0.56
1:B:110:ASP:OD1	1:B:264:HIS:HE1	1.89	0.54
1:D:110:ASP:OD1	1:D:264:HIS:HE1	1.91	0.54
1:D:58:ARG:CD	1:D:253:ASP:OD1	2.56	0.54
1:D:147:LEU:HB2	1:D:152:ARG:HG2	1.90	0.53
1:A:69:GLU:OE2	1:A:90:THR:HB	2.08	0.53
1:D:104:ALA:HB1	3:D:2025:HOH:O	2.10	0.51
1:C:163:SER:HB2	1:C:193:GLU:HG2	1.93	0.51
1:D:23:ARG:NH1	3:D:2002:HOH:O	2.44	0.50
1:C:110:ASP:OD1	1:C:264:HIS:CE1	2.64	0.50
1:C:33:LEU:HD22	1:C:71:PHE:CE1	2.48	0.48
1:C:279:ARG:HD3	1:C:281:THR:OG1	2.13	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:269:ASP:O	1:C:22:GLY:HA3	2.14	0.48
1:B:34:SER:HB2	1:B:35:PRO:CD	2.46	0.46
1:D:117:TYR:HB2	1:D:118:PRO:CD	2.46	0.46
1:B:60:GLN:HA	1:B:62:HIS:CE1	2.51	0.45
1:A:110:ASP:OD1	1:A:264:HIS:CE1	2.65	0.45
1:B:95:GLN:NE2	1:B:97:ASP:OD2	2.38	0.44
1:D:23:ARG:HH21	1:D:23:ARG:HD3	1.48	0.44
1:A:154:LEU:HD22	1:B:154:LEU:HD22	2.00	0.43
1:B:279:ARG:HD3	1:B:281:THR:OG1	2.19	0.43
1:C:78:PRO:HA	1:C:281:THR:O	2.19	0.42
1:A:153:GLU:OE1	1:B:157:LYS:NZ	2.26	0.42
1:D:117:TYR:HB2	1:D:118:PRO:HD2	2.02	0.42
1:B:110:ASP:OD1	1:B:264:HIS:CE1	2.71	0.41
1:B:69:GLU:OE2	1:B:90:THR:HB	2.20	0.41
1:B:285:GLU:H	1:B:285:GLU:CD	2.29	0.41
1:B:78:PRO:HA	1:B:281:THR:O	2.21	0.41
1:C:153:GLU:O	1:C:157:LYS:HD3	2.21	0.41
1:B:22:GLY:HA3	1:C:269:ASP:O	2.21	0.40
1:C:112:THR:O	1:C:196:HIS:NE2	2.54	0.40
1:B:299:ARG:HE	1:B:299:ARG:HB3	1.78	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	246/301 (82%)	239 (97%)	7 (3%)	0	100	100
1	B	246/301 (82%)	238 (97%)	8 (3%)	0	100	100
1	C	236/301 (78%)	227 (96%)	9 (4%)	0	100	100
1	D	236/301 (78%)	229 (97%)	7 (3%)	0	100	100
All	All	964/1204 (80%)	933 (97%)	31 (3%)	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	199/244 (82%)	194 (98%)	5 (2%)	42	45
1	B	197/244 (81%)	192 (98%)	5 (2%)	42	45
1	C	180/244 (74%)	172 (96%)	8 (4%)	25	23
1	D	179/244 (73%)	174 (97%)	5 (3%)	38	40
All	All	755/976 (77%)	732 (97%)	23 (3%)	36	38

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	ARG
1	A	96	LEU
1	A	111	VAL
1	A	124	ARG
1	A	214	HIS
1	B	79	VAL
1	B	96	LEU
1	B	111	VAL
1	B	124	ARG
1	B	191	VAL
1	C	61	THR
1	C	79	VAL
1	C	103	ARG
1	C	111	VAL
1	C	214	HIS
1	C	219	ILE
1	C	228	GLN
1	C	279	ARG
1	D	90	THR
1	D	96	LEU
1	D	111	VAL
1	D	149	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	214	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	HIS
1	A	264	HIS
1	B	264	HIS
1	D	145	GLN
1	D	264	HIS

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	SIN	B	1301	-	7,7,7	1.12	0	8,8,8	1.95	2 (25%)
2	SIN	D	1300	-	7,7,7	1.36	0	8,8,8	1.99	4 (50%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SIN	C	1300	-	7,7,7	1.36	2 (28%)	8,8,8	1.93	1 (12%)
2	SIN	A	1301	-	7,7,7	1.04	0	8,8,8	2.09	4 (50%)

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	SIN	B	1301	-	-	2/5/5/5	-
2	SIN	D	1300	-	-	2/5/5/5	-
2	SIN	C	1300	-	-	0/5/5/5	-
2	SIN	A	1301	-	-	2/5/5/5	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1300	SIN	O4-C4	-2.22	1.23	1.30
2	C	1300	SIN	O1-C1	2.14	1.29	1.22

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1300	SIN	C3-C2-C1	-4.36	102.09	113.67
2	B	1301	SIN	O4-C4-C3	3.34	124.55	114.00
2	A	1301	SIN	O4-C4-C3	3.30	124.42	114.00
2	D	1300	SIN	O4-C4-C3	3.07	123.69	114.00
2	D	1300	SIN	O4-C4-O3	-2.63	116.57	123.33
2	A	1301	SIN	O2-C1-C2	2.56	122.08	114.00
2	B	1301	SIN	O4-C4-O3	-2.38	117.22	123.33
2	A	1301	SIN	O4-C4-O3	-2.33	117.33	123.33
2	D	1300	SIN	C3-C2-C1	-2.32	107.50	113.67
2	A	1301	SIN	O2-C1-O1	-2.17	117.74	123.33
2	D	1300	SIN	O2-C1-O1	-2.13	117.85	123.33

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1301	SIN	C2-C3-C4-O4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
2	D	1300	SIN	O2-C1-C2-C3
2	D	1300	SIN	O1-C1-C2-C3
2	B	1301	SIN	C2-C3-C4-O3
2	A	1301	SIN	C2-C3-C4-O4
2	A	1301	SIN	C2-C3-C4-O3

There are no ring outliers.

No monomer is involved in short contacts.

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	254/301 (84%)	0.06	12 (4%) 36 41	14, 24, 44, 67	0
1	B	254/301 (84%)	0.03	10 (3%) 43 48	15, 24, 44, 63	0
1	C	244/301 (81%)	0.17	17 (6%) 22 25	14, 25, 46, 58	0
1	D	244/301 (81%)	0.56	30 (12%) 8 9	14, 26, 43, 56	0
All	All	996/1204 (82%)	0.20	69 (6%) 23 26	14, 25, 44, 67	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	80	ALA	7.7
1	D	192	TYR	6.6
1	D	299	ARG	6.5
1	B	80	ALA	6.0
1	A	300	LYS	5.9
1	A	191	VAL	5.8
1	C	96	LEU	5.5
1	C	192	TYR	5.3
1	D	61	THR	5.2
1	D	97	ASP	5.2
1	C	92	TYR	5.0
1	A	85	PRO	5.0
1	C	94	LEU	4.9
1	C	93	LEU	4.7
1	B	191	VAL	4.7
1	B	300	LYS	4.6
1	D	224	LEU	4.6
1	B	103	ARG	4.4
1	C	299	ARG	4.4
1	B	165	GLU	4.1
1	C	91	ARG	4.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	85	PRO	4.0
1	C	13	LEU	3.9
1	D	221	GLY	3.9
1	D	62	HIS	3.8
1	A	62	HIS	3.8
1	D	79	VAL	3.8
1	A	165	GLU	3.5
1	B	62	HIS	3.5
1	C	95	GLN	3.5
1	D	92	TYR	3.4
1	D	296	ARG	3.4
1	A	103	ARG	3.4
1	D	90	THR	3.3
1	D	298	THR	3.2
1	A	192	TYR	3.2
1	C	103	ARG	3.2
1	D	164	ASN	3.2
1	A	13	LEU	3.1
1	D	13	LEU	3.0
1	D	63	LEU	3.0
1	C	62	HIS	2.9
1	B	192	TYR	2.9
1	D	14	GLU	2.7
1	D	219	ILE	2.7
1	D	104	ALA	2.6
1	A	299	ARG	2.6
1	D	145	GLN	2.6
1	B	224	LEU	2.6
1	B	13	LEU	2.5
1	C	14	GLU	2.5
1	C	296	ARG	2.5
1	D	65	ASP	2.5
1	C	165	GLU	2.5
1	A	211	GLN	2.4
1	A	79	VAL	2.4
1	D	297	THR	2.3
1	D	94	LEU	2.3
1	D	211	GLN	2.2
1	D	67	SER	2.2
1	D	218	ARG	2.2
1	C	211	GLN	2.1
1	C	61	THR	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	224	LEU	2.1
1	D	161	VAL	2.1
1	D	228	GLN	2.0
1	D	66	GLN	2.0
1	D	230	LEU	2.0
1	D	64	ASP	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($\text{\AA}^2$)	Q<0.9
2	SIN	C	1300	8/8	0.90	0.14	29,37,54,56	0
2	SIN	D	1300	8/8	0.90	0.13	30,33,54,56	0
2	SIN	B	1301	8/8	0.92	0.12	32,34,42,45	0
2	SIN	A	1301	8/8	0.94	0.13	22,29,43,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.