



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:30 AM UTC

PDB ID : 7SME / pdb_00007sme
Title : p107 pocket domain complexed with HDAC1 peptide
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Deposited on : 2021-10-25
Resolution : 2.64 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

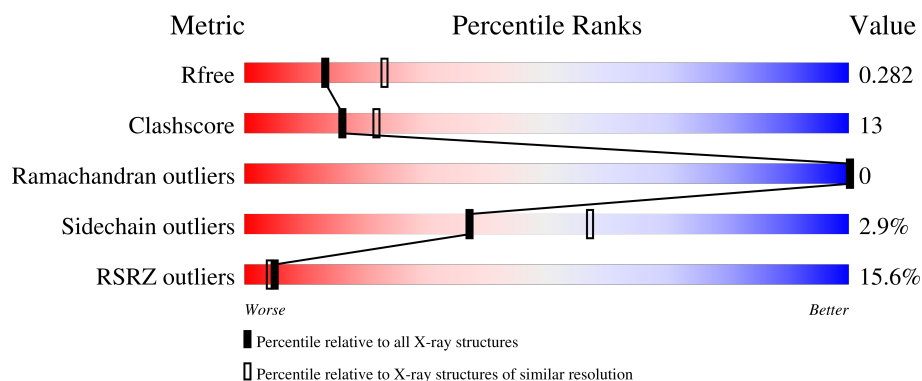
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.64 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	2053 (2.66-2.62)
Clashscore	190562	2097 (2.66-2.62)
Ramachandran outliers	187476	2066 (2.66-2.62)
Sidechain outliers	187428	2066 (2.66-2.62)
RSRZ outliers	180081	2052 (2.66-2.62)

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	371	
2	B	10	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2883 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 Retinoblastoma-like protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	342	2774	1785	466	505	18	0	0	0

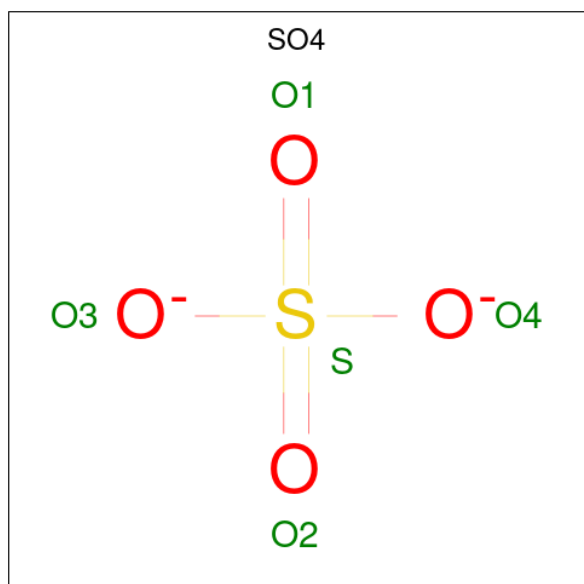
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	388	GLY	-	expression tag	UNP P28749
A	389	GLU	-	expression tag	UNP P28749
A	390	PHE	-	expression tag	UNP P28749

- Molecule 2 is a protein called Histone deacetylase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	10	68	42	10	15	1	0	0	0

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

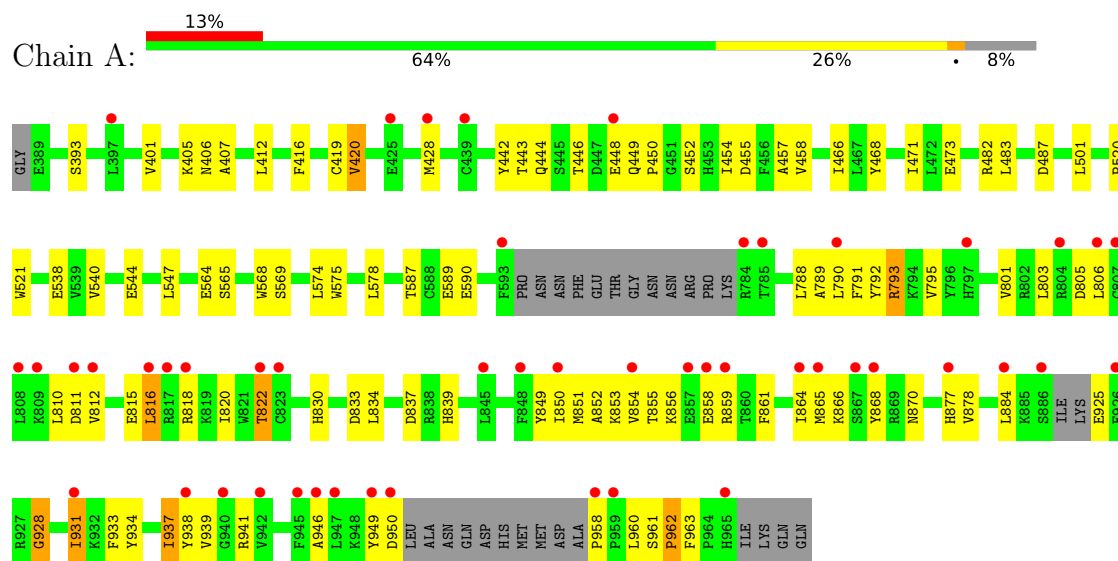
- Molecule 4 is water.

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

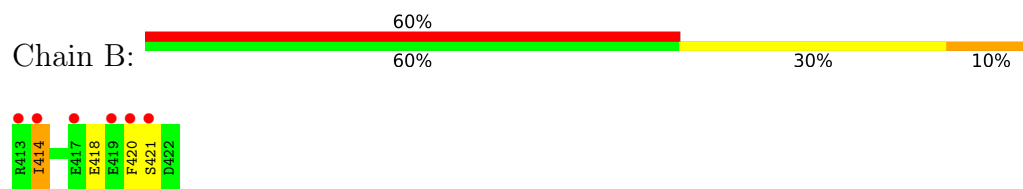
3 Residue-property plots [i](#)

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: Retinoblastoma-like protein 1



• Molecule 2: Histone deacetylase 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	99.91Å 75.17Å 71.74Å 90.00° 109.86° 90.00°	Depositor
Resolution (Å)	46.98 – 2.64 46.98 – 2.64	Depositor EDS
% Data completeness (in resolution range)	99.5 (46.98-2.64) 99.7 (46.98-2.64)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.29 (at 2.65Å)	Xtriage
Refinement program	PHENIX (1.17.1_3660: ???)	Depositor
R, R_{free}	0.224 , 0.277 0.229 , 0.282	Depositor DCC
R_{free} test set	709 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	47.5	Xtriage
Anisotropy	0.564	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 45.4	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	0.91	EDS
Total number of atoms	2883	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.92% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.99	0/2834	1.48	33/3837 (0.9%)
2	B	0.87	0/68	1.23	0/91
All	All	0.99	0/2902	1.47	33/3928 (0.8%)

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	443	THR	CA-C-N	-7.03	110.68	122.64
1	A	443	THR	C-N-CA	-7.03	110.68	122.64
1	A	830	HIS	CA-C-N	-6.48	115.72	122.85
1	A	830	HIS	C-N-CA	-6.48	115.72	122.85
1	A	568	TRP	CA-C-N	6.43	130.27	120.82
1	A	568	TRP	C-N-CA	6.43	130.27	120.82
1	A	928	GLY	CA-C-N	6.26	132.57	121.75
1	A	928	GLY	C-N-CA	6.26	132.57	121.75
1	A	406	ASN	N-CA-C	-6.22	105.96	112.93
1	A	565	SER	N-CA-C	-6.09	99.57	107.73
1	A	793	ARG	O-C-N	6.07	128.57	122.08
1	A	816	LEU	N-CA-C	-6.05	104.80	111.82
1	A	420	VAL	O-C-N	6.00	127.79	121.91
1	A	822	THR	CB-CA-C	5.88	120.85	110.85
1	A	565	SER	CA-C-O	-5.72	114.99	120.94
1	A	937	ILE	O-C-N	5.60	126.96	121.98
1	A	448	GLU	O-C-N	5.56	128.03	122.08
1	A	815	GLU	N-CA-C	-5.54	106.87	113.97
1	A	444	GLN	O-C-N	5.50	129.43	123.10
1	A	789	ALA	CA-C-N	-5.49	111.53	121.14
1	A	789	ALA	C-N-CA	-5.49	111.53	121.14
1	A	859	ARG	CA-C-N	-5.31	113.52	123.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	859	ARG	C-N-CA	-5.31	113.52	123.03
1	A	540	VAL	CA-C-O	-5.27	115.79	121.27
1	A	962	PRO	CA-C-N	5.26	128.92	121.61
1	A	962	PRO	C-N-CA	5.26	128.92	121.61
1	A	468	TYR	N-CA-C	-5.19	105.52	111.07
1	A	455	ASP	CB-CA-C	5.15	119.34	110.79
1	A	792	TYR	CB-CA-C	5.15	119.60	110.85
1	A	837	ASP	CB-CA-C	5.11	119.58	111.51
1	A	938	TYR	CA-C-N	-5.04	113.26	120.42
1	A	938	TYR	C-N-CA	-5.04	113.26	120.42
1	A	393	SER	O-C-N	5.02	127.25	122.03

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	2774	0	2747	74	0
2	B	68	0	50	2	0
3	A	10	0	0	0	0
4	A	31	0	0	2	0
All	All	2883	0	2797	74	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 13.

All (74) 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:790:LEU:HD12	1:A:793:ARG:HD3	1.48	0.93
1:A:446:THR:HG23	1:A:449:GLN:H	1.38	0.88
1:A:849:TYR:CE2	1:A:853:LYS:HD2	2.07	0.88
1:A:822:THR:HG22	1:A:961:SER:H	1.41	0.85
1:A:822:THR:CG2	1:A:961:SER:H	1.94	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:MET:O	1:A:428:MET:HE3	1.80	0.80
1:A:791:PHE:O	1:A:795:VAL:HG23	1.85	0.76
1:A:818:ARG:O	1:A:822:THR:HG23	1.84	0.75
1:A:589:GLU:HG3	1:A:962:PRO:CG	2.17	0.73
1:A:544:GLU:HG3	1:A:547:LEU:HD13	1.75	0.69
1:A:574:LEU:HD12	1:A:788:LEU:CD2	2.22	0.68
1:A:587:THR:OG1	1:A:590:GLU:HG3	1.94	0.67
1:A:574:LEU:HD12	1:A:788:LEU:HD21	1.76	0.67
1:A:589:GLU:HG3	1:A:962:PRO:HG3	1.78	0.66
1:A:446:THR:CG2	1:A:449:GLN:H	2.08	0.65
1:A:949:TYR:CE2	1:A:960:LEU:HD21	2.32	0.65
1:A:854:VAL:HB	1:A:950:ASP:OD1	1.97	0.64
1:A:442:TYR:CE1	1:A:452:SER:HB2	2.32	0.64
1:A:816:LEU:HG	1:A:820:ILE:HD11	1.78	0.64
1:A:852:ALA:HA	1:A:855:THR:OG1	1.98	0.64
1:A:939:VAL:HG11	2:B:414:ILE:O	1.99	0.62
1:A:820:ILE:HG13	1:A:851:MET:HB2	1.82	0.61
1:A:834:LEU:HD21	1:A:937:ILE:HG22	1.82	0.61
1:A:931:ILE:HD13	2:B:418:GLU:O	2.01	0.61
1:A:849:TYR:CZ	1:A:853:LYS:HD2	2.36	0.60
1:A:816:LEU:HD13	1:A:855:THR:HG21	1.83	0.60
1:A:454:ILE:O	1:A:458:VAL:HG23	2.02	0.59
1:A:816:LEU:O	1:A:820:ILE:HD12	2.01	0.59
1:A:482:ARG:NH2	1:A:483:LEU:HD23	2.18	0.58
1:A:884:LEU:HA	1:A:937:ILE:HD11	1.87	0.55
1:A:790:LEU:CD1	1:A:793:ARG:HD3	2.31	0.54
1:A:850:ILE:O	1:A:854:VAL:HG13	2.08	0.54
1:A:574:LEU:CD1	1:A:788:LEU:CD2	2.85	0.53
1:A:420:VAL:HG23	1:A:487:ASP:OD2	2.09	0.53
1:A:401:VAL:HB	1:A:466:ILE:HG21	1.92	0.51
1:A:811:ASP:O	1:A:811:ASP:OD1	2.29	0.51
1:A:934:TYR:O	1:A:939:VAL:HG23	2.11	0.51
1:A:578:LEU:HA	1:A:788:LEU:HD11	1.92	0.50
1:A:482:ARG:NH2	1:A:483:LEU:CD2	2.75	0.49
1:A:833:ASP:OD2	1:A:941:ARG:NH1	2.46	0.49
1:A:849:TYR:CG	1:A:861:PHE:CE1	3.00	0.48
1:A:544:GLU:CG	1:A:547:LEU:HD13	2.43	0.48
1:A:442:TYR:CD1	1:A:457:ALA:HB2	2.49	0.46
1:A:810:LEU:HD11	1:A:864:ILE:HA	1.97	0.46
1:A:851:MET:HA	1:A:854:VAL:HG22	1.97	0.46
1:A:816:LEU:HG	1:A:820:ILE:CD1	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:884:LEU:HD11	1:A:928:GLY:N	2.30	0.46
1:A:520:PRO:O	1:A:521:TRP:C	2.59	0.46
1:A:428:MET:HE1	4:A:1111:HOH:O	2.16	0.46
1:A:949:TYR:CE2	1:A:958:PRO:HB2	2.51	0.46
1:A:569:SER:O	1:A:575:TRP:NE1	2.48	0.45
1:A:790:LEU:HA	1:A:793:ARG:HB2	1.98	0.45
1:A:849:TYR:OH	1:A:853:LYS:CE	2.65	0.45
1:A:933:PHE:CD1	1:A:937:ILE:HD12	2.51	0.45
1:A:820:ILE:HD12	1:A:820:ILE:H	1.82	0.45
1:A:868:TYR:C	1:A:870:ASN:N	2.74	0.45
1:A:405:LYS:C	1:A:407:ALA:H	2.26	0.44
1:A:564:GLU:O	1:A:839:HIS:HA	2.17	0.44
1:A:820:ILE:HG13	1:A:851:MET:CB	2.46	0.44
1:A:856:LYS:C	1:A:858:GLU:OE1	2.60	0.44
1:A:412:LEU:HG	1:A:416:PHE:CZ	2.54	0.43
1:A:806:LEU:HA	1:A:806:LEU:HD23	1.80	0.43
1:A:854:VAL:HG11	1:A:946:ALA:HB1	2.00	0.43
1:A:801:VAL:O	1:A:805:ASP:OD2	2.37	0.43
1:A:933:PHE:HD1	1:A:937:ILE:HD12	1.83	0.43
1:A:412:LEU:HD13	1:A:473:GLU:HA	2.01	0.42
1:A:471:ILE:HG21	1:A:501:LEU:HD11	2.01	0.42
1:A:450:PRO:HG2	4:A:1109:HOH:O	2.20	0.42
1:A:822:THR:HG21	1:A:961:SER:H	1.77	0.41
1:A:803:LEU:HD23	1:A:963:PHE:HE1	1.85	0.41
1:A:856:LYS:O	1:A:858:GLU:OE1	2.39	0.41
1:A:419:CYS:HA	1:A:487:ASP:OD2	2.21	0.41
1:A:877:HIS:CE1	1:A:878:VAL:HG23	2.55	0.40
1:A:834:LEU:HD21	1:A:937:ILE:CG2	2.49	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	334/371 (90%)	323 (97%)	11 (3%)	0	100	100
2	B	8/10 (80%)	6 (75%)	2 (25%)	0	100	100
All	All	342/381 (90%)	329 (96%)	13 (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	309/342 (90%)	303 (98%)	6 (2%)	50	70
2	B	6/9 (67%)	3 (50%)	3 (50%)	0	0
All	All	315/351 (90%)	306 (97%)	9 (3%)	37	57

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	538	GLU
1	A	812	VAL
1	A	865	MET
1	A	866	LYS
1	A	925	GLU
1	A	931	ILE
2	B	414	ILE
2	B	420	PHE
2	B	421	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	398	GLN
1	A	459	ASN
1	A	871	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](#)

2 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	$\# Z > 2$
3	SO4	A	1002	-	4,4,4	0.31	0	6,6,6	0.28	0
3	SO4	A	1001	-	4,4,4	0.28	0	6,6,6	0.23	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.

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 ⓘ

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	342/371 (92%)	0.84	49 (14%) 6 5	30, 50, 93, 100	0
2	B	10/10 (100%)	2.61	6 (60%) 0 0	95, 100, 102, 106	0
All	All	352/381 (92%)	0.89	55 (15%) 5 4	30, 50, 96, 106	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	414	ILE	4.4
1	A	965	HIS	4.2
2	B	419	GLU	4.0
1	A	854	VAL	4.0
1	A	848	PHE	3.8
2	B	413	ARG	3.6
1	A	886	SER	3.6
2	B	421	SER	3.5
1	A	857	GLU	3.5
1	A	784	ARG	3.5
1	A	593	PHE	3.4
1	A	864	ILE	3.3
1	A	428	MET	3.2
1	A	806	LEU	3.2
1	A	947	LEU	3.2
1	A	958	PRO	3.2
1	A	940	GLY	3.1
1	A	946	ALA	3.0
1	A	959	PRO	3.0
1	A	818	ARG	3.0
1	A	865	MET	2.9
1	A	950	ASP	2.8
1	A	845	LEU	2.7
1	A	859	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	420	PHE	2.6
1	A	850	ILE	2.6
1	A	931	ILE	2.6
1	A	809	LYS	2.5
1	A	397	LEU	2.5
1	A	942	VAL	2.4
1	A	812	VAL	2.4
2	B	417	GLU	2.4
1	A	439	CYS	2.4
1	A	425	GLU	2.3
1	A	808	LEU	2.3
1	A	823	CYS	2.2
1	A	858	GLU	2.2
1	A	868	TYR	2.2
1	A	938	TYR	2.2
1	A	867	SER	2.2
1	A	817	ARG	2.2
1	A	926	GLU	2.2
1	A	807	CYS	2.2
1	A	790	LEU	2.2
1	A	945	PHE	2.2
1	A	785	THR	2.1
1	A	797	HIS	2.1
1	A	949	TYR	2.1
1	A	811	ASP	2.1
1	A	448	GLU	2.0
1	A	822	THR	2.0
1	A	816	LEU	2.0
1	A	884	LEU	2.0
1	A	877	HIS	2.0
1	A	804	ARG	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
3	SO4	A	1001	5/5	0.78	0.18	72,74,91,91	0
3	SO4	A	1002	5/5	0.98	0.06	32,32,43,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.