



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:46 AM UTC

PDB ID : 7S2L / pdb_00007s2l
Title : Crystal structure of sulfonamide resistance enzyme Sul3 apoenzyme
Authors : Stogios, P.J.; Venkatesan, M.; Michalska, K.; Mesa, N.; Di Leo, R.; Savchenko, A.; Joachimiak, A.; Satchell, K.J.F.; Center for Structural Biology of Infectious Diseases (CSBID); Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2021-09-03
Resolution : 2.79 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

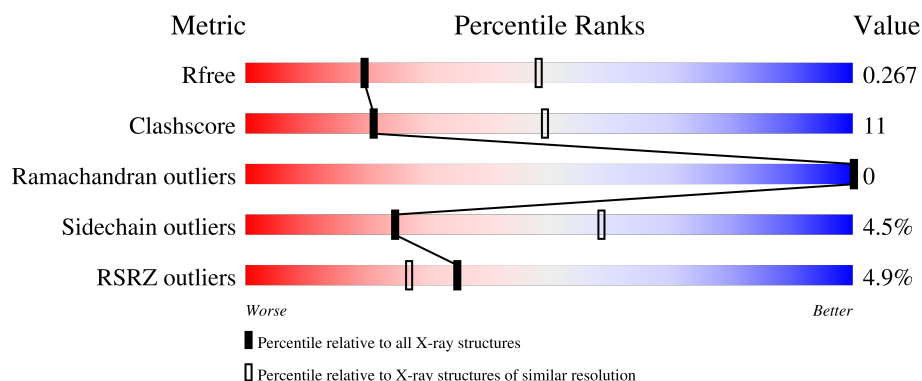
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	265	5% 72% 27% .
1	B	265	5% 67% 29% ..
1	C	265	5% 77% 22% .
1	D	265	5% 68% 25% 6%

2 Entry composition [i](#)

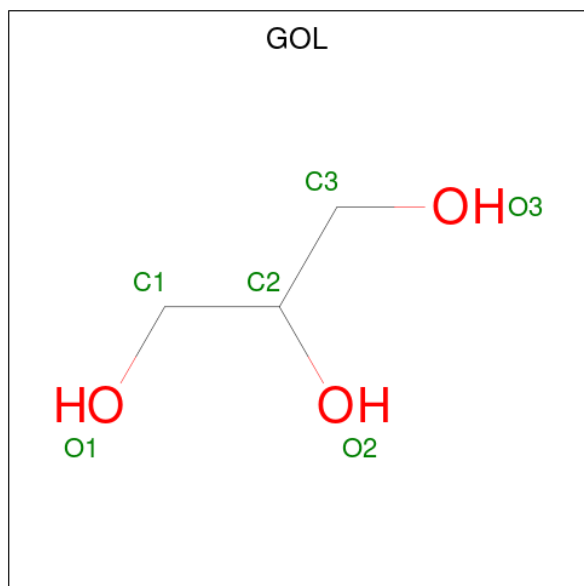
There are 5 unique types of molecules in this entry. The entry contains 8184 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Sul3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	0	0
			2032	1299	334	390	9			
1	B	260	Total	C	N	O	S	0	0	0
			1992	1275	328	380	9			
1	C	264	Total	C	N	O	S	0	1	0
			2029	1298	332	390	9			
1	D	249	Total	C	N	O	S	0	0	0
			1912	1224	316	363	9			

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		

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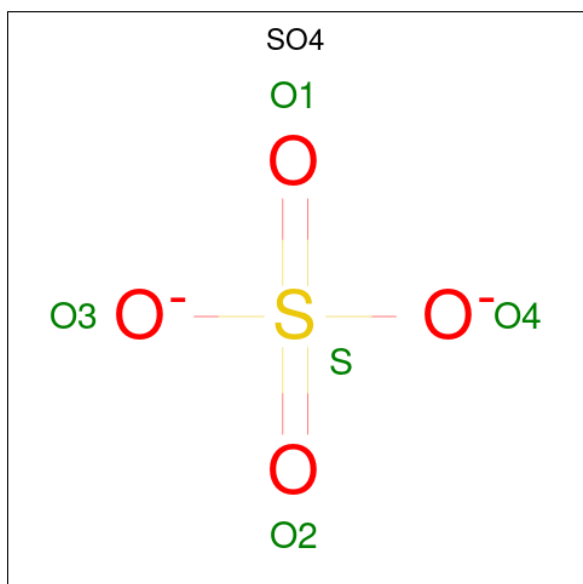
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	C	2	Total	Cl	0	0
			2	2		
3	D	1	Total	Cl	0	0
			1	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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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	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	C	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		

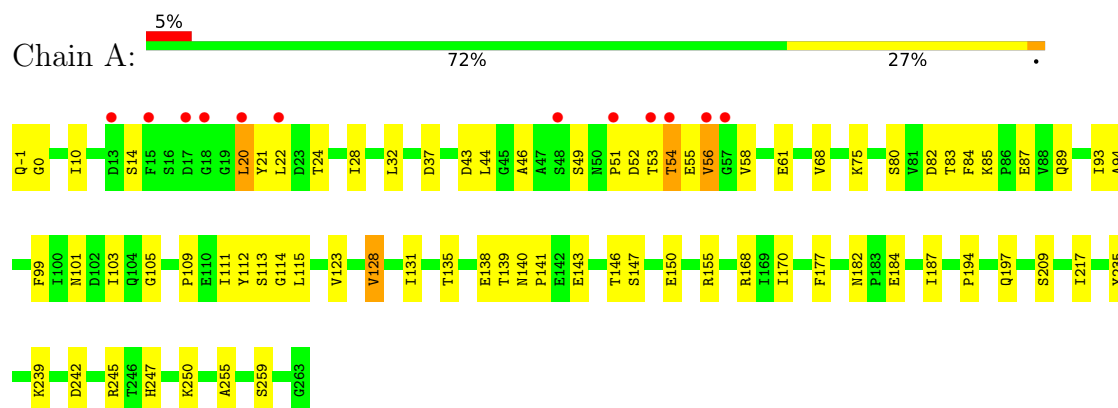
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	37	Total	O	0	0
			37	37		
5	B	31	Total	O	0	0
			31	31		
5	C	37	Total	O	0	0
			37	37		
5	D	23	Total	O	0	0
			23	23		

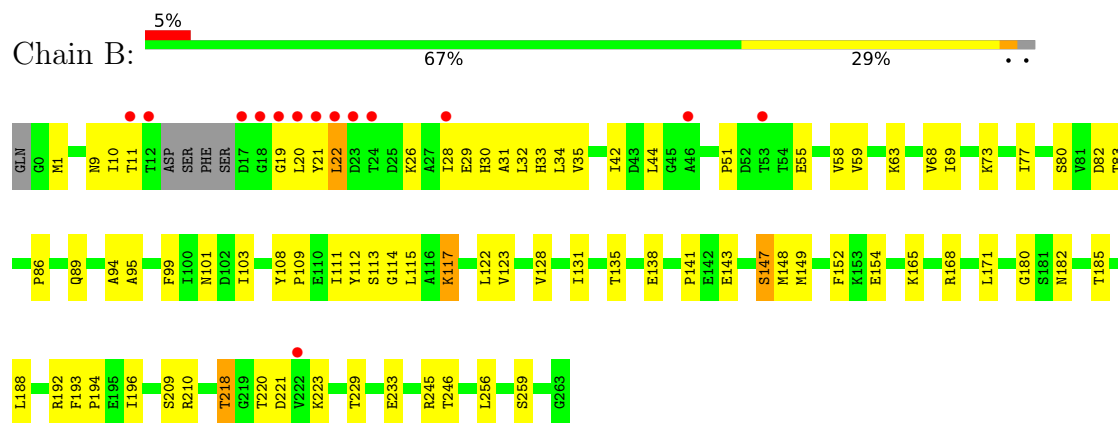
3 Residue-property plots

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.

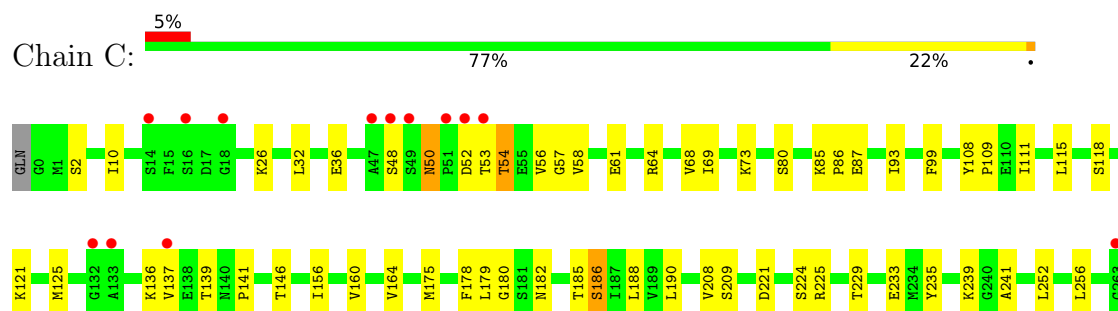
• Molecule 1: Sul3



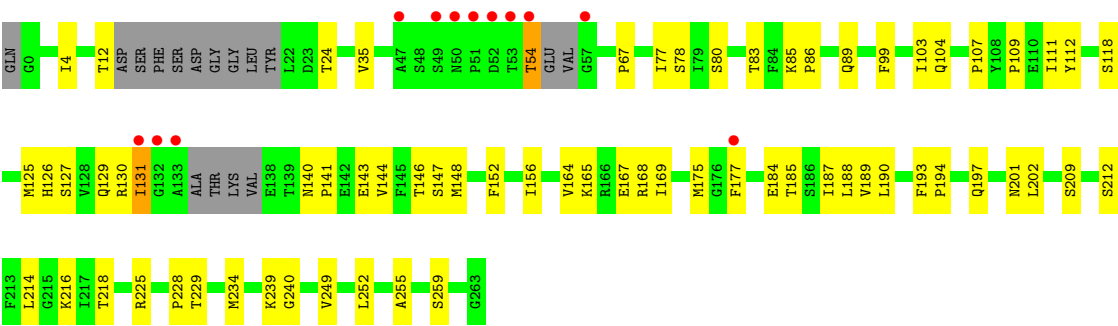
• Molecule 1: Sul3



• Molecule 1: Sul3



● Molecule 1: Sul3



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	123.80Å 123.80Å 434.42Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.79 30.00 – 2.79	Depositor EDS
% Data completeness (in resolution range)	99.4 (30.00-2.79) 88.9 (30.00-2.79)	Depositor EDS
R_{merge}	0.31	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.73 (at 2.80Å)	Xtriage
Refinement program	phenix.refine 1.19.2_4158, PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.226 , 0.270 0.226 , 0.267	Depositor DCC
R_{free} test set	2000 reflections (4.02%)	wwPDB-VP
Wilson B-factor (Å ²)	55.8	Xtriage
Anisotropy	0.610	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8184	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.42% 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: SO4, CL, GOL

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	0.23	0/2065	0.44	0/2784
1	B	0.22	0/2023	0.42	0/2726
1	C	0.25	0/2065	0.43	0/2784
1	D	0.22	0/1940	0.40	0/2610
All	All	0.23	0/8093	0.42	0/10904

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	2032	0	2081	50	0
1	B	1992	0	2049	54	0
1	C	2029	0	2079	47	0
1	D	1912	0	1968	41	0
2	A	18	0	24	0	0
2	C	18	0	24	0	0
2	D	6	0	8	1	0
3	A	1	0	0	0	0
3	C	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	1	0	0	0	0
4	A	20	0	0	1	0
4	B	10	0	0	1	0
4	C	10	0	0	0	0
4	D	5	0	0	0	0
5	A	37	0	0	1	0
5	B	31	0	0	1	0
5	C	37	0	0	1	0
5	D	23	0	0	0	0
All	All	8184	0	8233	179	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 11.

All (179) 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:149:MET:HE1	1:B:196:ILE:HG12	1.50	0.93
1:C:179:LEU:HB3	1:C:185:THR:HG21	1.57	0.85
1:D:218:THR:HG21	1:D:228:PRO:HB3	1.61	0.82
1:A:54:THR:HG21	1:C:54:THR:HB	1.63	0.81
1:B:103:ILE:HD11	1:B:131:ILE:HD11	1.64	0.78
1:C:53:THR:HG23	1:C:85:LYS:HD3	1.67	0.77
1:B:218:THR:HG22	1:B:220:THR:H	1.51	0.75
1:D:54:THR:HA	1:D:85:LYS:HE2	1.66	0.75
1:C:179:LEU:HB3	1:C:185:THR:CG2	2.17	0.75
1:C:141:PRO:HG3	1:C:182:ASN:HD21	1.51	0.74
1:B:245:ARG:NH2	4:B:302:SO4:O3	2.22	0.73
1:A:245:ARG:NH2	4:A:307:SO4:O3	2.23	0.72
1:A:56:VAL:O	1:A:85:LYS:HE2	1.90	0.70
1:C:50:ASN:N	1:C:50:ASN:OD1	2.25	0.70
1:B:188:LEU:HD11	1:B:192:ARG:HE	1.56	0.69
1:D:104:GLN:HB3	1:D:107:PRO:HG3	1.75	0.69
1:A:37:ASP:HB3	1:A:250:LYS:HB2	1.75	0.68
1:B:220:THR:HG22	1:B:221:ASP:H	1.58	0.67
1:A:54:THR:CG2	1:C:54:THR:HB	2.25	0.66
1:D:86:PRO:HB3	1:D:111:ILE:HG12	1.77	0.66
1:A:82:ASP:HA	1:A:101:ASN:HB3	1.78	0.66
1:A:43:ASP:OD2	1:A:245:ARG:NH1	2.30	0.64
1:C:235:TYR:OH	1:C:239:LYS:NZ	2.29	0.64
1:C:179:LEU:HD13	1:C:185:THR:HG22	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:ILE:HG13	1:B:94:ALA:HB1	1.81	0.63
1:B:141:PRO:HA	1:B:185:THR:HG22	1.79	0.63
1:C:58:VAL:HG22	1:C:85:LYS:HD2	1.81	0.62
1:C:141:PRO:HG3	1:C:182:ASN:ND2	2.14	0.62
1:B:51:PRO:HB3	1:B:55:GLU:HG2	1.82	0.62
1:B:10:ILE:HA	1:B:22:LEU:HB2	1.82	0.60
1:A:20:LEU:O	1:A:20:LEU:HG	2.00	0.60
1:D:165:LYS:HE2	1:D:167:GLU:HB2	1.84	0.59
1:A:44:LEU:HD11	1:A:68:VAL:HG11	1.85	0.58
1:A:58:VAL:HG11	1:A:87:GLU:HB2	1.86	0.58
1:C:179:LEU:CD1	1:C:185:THR:HG22	2.33	0.58
1:A:84:PHE:CD1	1:A:103:ILE:HD11	2.38	0.58
1:A:146:THR:O	1:A:150:GLU:HG3	2.04	0.58
1:D:54:THR:O	1:D:54:THR:OG1	2.22	0.57
1:B:44:LEU:HD21	1:B:68:VAL:HG11	1.87	0.57
1:A:168:ARG:NH2	5:A:401:HOH:O	2.38	0.57
1:C:80:SER:HB2	1:C:99:PHE:HB2	1.87	0.56
1:D:240:GLY:HA2	2:D:301:GOL:H12	1.88	0.55
1:D:141:PRO:HA	1:D:185:THR:HG22	1.88	0.54
1:D:193:PHE:CD2	1:D:194:PRO:HD3	2.42	0.54
1:C:125:MET:HE1	1:C:178:PHE:CD2	2.42	0.54
1:A:209:SER:OG	1:A:247:HIS:ND1	2.37	0.54
1:D:118:SER:O	1:D:168:ARG:NH1	2.41	0.53
1:B:29:GLU:HA	1:B:32:LEU:HB2	1.90	0.53
1:D:175:MET:SD	1:D:190:LEU:HG	2.49	0.53
1:B:223:LYS:H	1:B:223:LYS:HD2	1.73	0.53
1:D:4:ILE:HG12	1:D:252:LEU:HD23	1.90	0.52
1:D:4:ILE:HG21	1:D:249:VAL:HG13	1.92	0.52
1:D:209:SER:HA	1:D:229:THR:HG23	1.92	0.52
1:A:80:SER:CB	1:A:99:PHE:HB2	2.40	0.52
1:D:156:ILE:H	1:D:156:ILE:HD12	1.74	0.52
1:A:24:THR:O	1:A:28:ILE:HG13	2.10	0.52
1:B:218:THR:HG22	1:B:220:THR:N	2.23	0.51
1:A:93:ILE:HD11	1:A:115:LEU:HD23	1.93	0.51
1:A:128:VAL:HG21	1:A:138:GLU:HG3	1.92	0.51
1:C:54:THR:HG23	1:C:57:GLY:H	1.75	0.51
1:B:31:ALA:HB1	1:B:42:ILE:HG12	1.94	0.50
1:A:140:ASN:HB3	1:A:143:GLU:HB2	1.93	0.50
1:A:54:THR:CG2	1:C:54:THR:CB	2.90	0.50
1:A:51:PRO:HG2	1:A:131:ILE:HG23	1.93	0.50
1:C:125:MET:HE1	1:C:178:PHE:CE2	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:128:VAL:HG21	1:B:138:GLU:HG3	1.94	0.49
1:A:52:ASP:HB2	1:C:56:VAL:HG21	1.93	0.49
1:D:234:MET:SD	1:D:255:ALA:HB1	2.52	0.49
1:B:69:ILE:O	1:B:73:LYS:HG2	2.12	0.49
1:D:175:MET:HE2	1:D:189:VAL:HG23	1.93	0.49
1:B:165:LYS:HD2	1:B:168:ARG:HD2	1.95	0.49
1:C:221:ASP:OD2	1:C:224:SER:OG	2.31	0.49
1:C:209:SER:HA	1:C:229:THR:HG23	1.94	0.49
1:C:221:ASP:O	1:C:225:ARG:HB2	2.13	0.49
1:B:28:ILE:HG13	1:B:68:VAL:HG13	1.95	0.48
1:C:69:ILE:O	1:C:73:LYS:HG2	2.14	0.48
1:D:83:THR:HG23	1:D:89:GLN:HG2	1.95	0.48
1:B:109:PRO:HA	1:B:112:TYR:CD1	2.48	0.48
1:D:184:GLU:HA	1:D:187:ILE:HD12	1.96	0.48
1:C:64:ARG:NH2	5:C:401:HOH:O	2.46	0.48
1:A:123:VAL:HG12	1:A:170:ILE:HB	1.95	0.48
1:B:9:ASN:HB3	1:B:11:THR:HG22	1.95	0.48
1:C:32:LEU:O	1:C:36:GLU:HG3	2.14	0.48
1:C:186:SER:O	1:C:190:LEU:HB2	2.13	0.48
1:B:26:LYS:O	1:B:30:HIS:N	2.40	0.48
1:B:109:PRO:HA	1:B:112:TYR:CG	2.49	0.48
1:D:109:PRO:HA	1:D:112:TYR:CD1	2.49	0.48
1:B:117:LYS:NZ	5:B:401:HOH:O	2.46	0.47
1:A:109:PRO:HA	1:A:112:TYR:CD1	2.48	0.47
1:B:210:ARG:HD3	1:B:223:LYS:O	2.14	0.47
1:D:35:VAL:HG21	1:D:77:ILE:HD13	1.97	0.47
1:B:83:THR:HG23	1:B:89:GLN:HG2	1.97	0.47
1:C:252:LEU:O	1:C:256:LEU:HD12	2.14	0.47
1:A:83:THR:HG23	1:A:89:GLN:HG2	1.97	0.46
1:D:24:THR:HG21	1:D:67:PRO:HG2	1.97	0.46
1:C:109:PRO:HG3	1:D:109:PRO:HD3	1.98	0.46
1:C:109:PRO:HG3	1:D:109:PRO:CD	2.45	0.46
1:A:235:TYR:CZ	1:A:239:LYS:HD2	2.51	0.46
1:A:10:ILE:HD11	1:A:44:LEU:HD12	1.97	0.46
1:B:80:SER:CB	1:B:99:PHE:HB2	2.46	0.46
1:A:255:ALA:O	1:A:259:SER:OG	2.27	0.45
1:D:103:ILE:HA	1:D:125:MET:CE	2.46	0.45
1:D:201:ASN:O	1:D:202:LEU:HD23	2.15	0.45
1:C:61:GLU:OE1	1:C:85:LYS:HE2	2.16	0.45
1:C:233:GLU:OE1	1:C:252:LEU:HD13	2.16	0.45
1:B:10:ILE:HG13	1:B:22:LEU:HD13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:136:LYS:HE2	1:C:137:VAL:HG22	1.97	0.45
1:B:29:GLU:O	1:B:33:HIS:ND1	2.49	0.45
1:C:180:GLY:C	1:C:182:ASN:H	2.24	0.45
1:B:19:GLY:C	1:B:20:LEU:HD22	2.42	0.45
1:C:50:ASN:O	1:C:52:ASP:N	2.50	0.45
1:B:82:ASP:HA	1:B:101:ASN:HB3	1.99	0.44
1:C:93:ILE:HD11	1:C:115:LEU:HD23	1.98	0.44
1:C:108:TYR:O	1:C:111:ILE:HG22	2.17	0.44
1:C:10:ILE:HD13	1:C:68:VAL:HG21	2.00	0.44
1:A:46:ALA:HB3	1:A:61:GLU:HG3	1.99	0.44
1:A:101:ASN:HA	1:A:123:VAL:HG22	1.98	0.44
1:A:114:GLY:HA3	1:B:94:ALA:HA	1.99	0.44
1:C:175:MET:HG2	1:C:208:VAL:HG23	1.99	0.44
1:C:93:ILE:HG23	1:C:118:SER:HB2	1.98	0.44
1:B:115:LEU:HD22	1:B:122:LEU:HD21	2.00	0.44
1:B:180:GLY:O	1:B:182:ASN:N	2.50	0.44
1:B:141:PRO:HB2	1:B:188:LEU:HD23	1.99	0.43
1:A:49:SER:O	1:A:51:PRO:HD3	2.19	0.43
1:A:85:LYS:O	1:A:89:GLN:HG3	2.18	0.43
1:B:35:VAL:HG21	1:B:77:ILE:HG21	1.99	0.43
1:D:140:ASN:O	1:D:144:VAL:HG23	2.18	0.43
1:A:105:GLY:O	1:A:155:ARG:NH1	2.42	0.43
1:A:94:ALA:HA	1:B:114:GLY:HA3	2.00	0.43
1:C:141:PRO:HB2	1:C:188:LEU:HD12	2.01	0.43
1:B:86:PRO:HG3	1:B:108:TYR:CD2	2.54	0.43
1:B:73:LYS:NZ	1:B:95:ALA:O	2.47	0.43
1:D:239:LYS:HD3	1:D:239:LYS:HA	1.74	0.43
1:D:126:HIS:CE1	1:D:148:MET:HB2	2.54	0.42
1:A:54:THR:HB	1:C:54:THR:OG1	2.18	0.42
1:A:80:SER:HB2	1:A:99:PHE:HB2	2.00	0.42
1:D:127:SER:OG	1:D:129:GLN:OE1	2.35	0.42
1:D:177:PHE:CE1	1:D:212:SER:HB2	2.54	0.42
1:A:32:LEU:HD21	1:A:75:LYS:HD2	2.01	0.42
1:B:30:HIS:CE1	1:B:34:LEU:HD11	2.55	0.42
1:C:99:PHE:CD2	1:C:121:LYS:HB2	2.54	0.42
1:D:216:LYS:HE3	1:D:216:LYS:HA	2.02	0.42
1:A:197:GLN:NE2	1:A:242:ASP:OD2	2.51	0.42
1:A:-1:GLN:HG2	1:A:0:GLY:H	1.84	0.42
1:A:14:SER:OG	1:A:21:TYR:O	2.38	0.42
1:B:209:SER:HA	1:B:229:THR:HG23	2.01	0.42
1:A:139:THR:O	1:A:141:PRO:HD3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:80:SER:CB	1:D:99:PHE:HB2	2.50	0.42
1:A:94:ALA:HB1	1:B:111:ILE:HG22	2.02	0.41
1:A:187:ILE:HD11	1:A:217:ILE:HD11	2.02	0.41
1:A:194:PRO:HG2	1:A:239:LYS:HE2	2.02	0.41
1:B:80:SER:HB2	1:B:99:PHE:HB2	2.02	0.41
1:A:135:THR:HG23	1:A:177:PHE:HE1	1.86	0.41
1:B:148:MET:SD	1:B:149:MET:HE2	2.60	0.41
1:C:80:SER:CB	1:C:99:PHE:HB2	2.48	0.41
1:D:152:PHE:O	1:D:156:ILE:HD12	2.20	0.41
1:A:37:ASP:CB	1:A:250:LYS:HB2	2.48	0.41
1:B:63:LYS:HD3	1:B:63:LYS:HA	1.92	0.41
1:B:143:GLU:O	1:B:147:SER:HB3	2.20	0.41
1:B:152:PHE:CD1	1:B:171:LEU:HD22	2.55	0.41
1:D:109:PRO:HA	1:D:112:TYR:CG	2.56	0.41
1:C:86:PRO:HG2	1:D:131:ILE:HG22	2.02	0.41
1:C:2:SER:OG	1:C:241:ALA:O	2.30	0.41
1:B:20:LEU:HD23	1:B:210:ARG:HH22	1.86	0.41
1:D:140:ASN:O	1:D:143:GLU:HG2	2.20	0.41
1:A:182:ASN:OD1	1:A:184:GLU:N	2.54	0.41
1:B:30:HIS:O	1:B:33:HIS:HB2	2.21	0.41
1:B:193:PHE:CD2	1:B:194:PRO:HD3	2.56	0.41
1:D:194:PRO:HA	1:D:197:GLN:OE1	2.21	0.41
1:D:214:LEU:O	1:D:218:THR:HG23	2.20	0.41
1:A:54:THR:HB	1:C:54:THR:HG1	1.86	0.40
1:B:1:MET:HE1	1:B:256:LEU:HB3	2.03	0.40
1:D:169:ILE:O	1:D:202:LEU:HD13	2.21	0.40
1:B:101:ASN:HA	1:B:123:VAL:HB	2.03	0.40
1:C:156:ILE:O	1:C:160:VAL:HG12	2.21	0.40
1:B:233:GLU:OE2	1:B:246:THR:OG1	2.25	0.40
1:C:58:VAL:HG11	1:C:87[B]:GLU:HB3	2.03	0.40
1:D:126:HIS:CD2	1:D:148:MET:HB2	2.56	0.40
1:B:193:PHE:CG	1:B:194:PRO:HD3	2.57	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	263/265 (99%)	253 (96%)	10 (4%)	0	100	100
1	B	256/265 (97%)	238 (93%)	18 (7%)	0	100	100
1	C	263/265 (99%)	251 (95%)	12 (5%)	0	100	100
1	D	241/265 (91%)	230 (95%)	11 (5%)	0	100	100
All	All	1023/1060 (96%)	972 (95%)	51 (5%)	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	222/222 (100%)	213 (96%)	9 (4%)	27	62
1	B	217/222 (98%)	206 (95%)	11 (5%)	21	54
1	C	222/222 (100%)	214 (96%)	8 (4%)	31	66
1	D	209/222 (94%)	198 (95%)	11 (5%)	20	52
All	All	870/888 (98%)	831 (96%)	39 (4%)	24	58

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
1	A	22	LEU
1	A	53	THR
1	A	54	THR
1	A	55	GLU
1	A	56	VAL
1	A	113	SER
1	A	128	VAL

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Mol	Chain	Res	Type
1	A	147	SER
1	B	21	TYR
1	B	22	LEU
1	B	58	VAL
1	B	59	VAL
1	B	113	SER
1	B	117	LYS
1	B	135	THR
1	B	147	SER
1	B	154	GLU
1	B	218	THR
1	B	259	SER
1	C	26	LYS
1	C	48	SER
1	C	50	ASN
1	C	54	THR
1	C	139	THR
1	C	146	THR
1	C	164	VAL
1	C	186	SER
1	D	12	THR
1	D	54	THR
1	D	78	SER
1	D	130	ARG
1	D	131	ILE
1	D	146	THR
1	D	147	SER
1	D	164	VAL
1	D	188	LEU
1	D	225	ARG
1	D	259	SER

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	89	GLN
1	B	30	HIS
1	C	182	ASN
1	D	30	HIS
1	D	89	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 20 ligands modelled in this entry, 4 are monoatomic - leaving 16 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	A	307	-	4,4,4	0.24	0	6,6,6	0.22	0
4	SO4	B	301	-	4,4,4	0.23	0	6,6,6	0.45	0
2	GOL	A	303	-	5,5,5	0.95	0	5,5,5	1.05	0
4	SO4	C	306	-	4,4,4	0.26	0	6,6,6	0.13	0
2	GOL	A	302	-	5,5,5	0.97	0	5,5,5	1.04	0
2	GOL	C	302	-	5,5,5	1.00	0	5,5,5	0.98	0
4	SO4	A	306	-	4,4,4	0.25	0	6,6,6	0.39	0
4	SO4	B	302	-	4,4,4	0.26	0	6,6,6	0.17	0
4	SO4	C	307	-	4,4,4	0.23	0	6,6,6	0.32	0
2	GOL	D	301	-	5,5,5	1.10	0	5,5,5	1.01	0
4	SO4	A	308	-	4,4,4	0.24	0	6,6,6	0.12	0
4	SO4	D	303	-	4,4,4	0.24	0	6,6,6	0.08	0
4	SO4	A	305	-	4,4,4	0.24	0	6,6,6	0.12	0
2	GOL	C	303	-	5,5,5	0.93	0	5,5,5	1.05	0
2	GOL	A	301	-	5,5,5	0.97	0	5,5,5	1.03	0
2	GOL	C	301	-	5,5,5	1.00	0	5,5,5	1.07	0

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	GOL	A	303	-	-	0/4/4/4	-
2	GOL	A	302	-	-	0/4/4/4	-
2	GOL	C	302	-	-	2/4/4/4	-
2	GOL	D	301	-	-	0/4/4/4	-
2	GOL	C	303	-	-	2/4/4/4	-
2	GOL	A	301	-	-	0/4/4/4	-
2	GOL	C	301	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	301	GOL	C1-C2-C3-O3
2	C	303	GOL	O1-C1-C2-C3
2	C	303	GOL	O1-C1-C2-O2
2	C	302	GOL	O1-C1-C2-C3
2	C	301	GOL	O2-C2-C3-O3
2	C	301	GOL	O1-C1-C2-O2
2	C	302	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	307	SO4	1	0
4	B	302	SO4	1	0
2	D	301	GOL	1	0

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	265/265 (100%)	0.08	12 (4%)	38 30	48, 63, 131, 262	0
1	B	260/265 (98%)	0.30	14 (5%)	31 24	50, 66, 117, 184	0
1	C	264/265 (99%)	0.15	13 (4%)	35 27	50, 68, 117, 210	1 (0%)
1	D	249/265 (93%)	0.25	12 (4%)	35 28	57, 78, 122, 225	0
All	All	1038/1060 (97%)	0.19	51 (4%)	35 27	48, 68, 124, 262	1 (0%)

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	19	GLY	7.2
1	A	56	VAL	7.0
1	B	21	TYR	5.4
1	B	20	LEU	5.3
1	D	47	ALA	5.2
1	A	20	LEU	5.1
1	C	14	SER	4.8
1	B	23	ASP	4.8
1	A	54	THR	4.3
1	D	131	ILE	4.1
1	A	22	LEU	3.8
1	D	54	THR	3.7
1	A	53	THR	3.4
1	D	57	GLY	3.4
1	B	11	THR	3.3
1	B	12	THR	3.3
1	C	51	PRO	3.2
1	D	51	PRO	3.2
1	D	132	GLY	3.1
1	D	49	SER	3.1
1	B	53	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	18	GLY	3.0
1	C	49	SER	2.9
1	D	52	ASP	2.9
1	D	177	PHE	2.9
1	C	47	ALA	2.8
1	A	57	GLY	2.8
1	B	17	ASP	2.8
1	A	51	PRO	2.8
1	D	53	THR	2.8
1	B	18	GLY	2.7
1	B	22	LEU	2.6
1	A	18	GLY	2.6
1	C	263	GLY	2.5
1	B	46	ALA	2.5
1	D	133	ALA	2.5
1	C	132	GLY	2.5
1	C	137	VAL	2.5
1	A	15	PHE	2.5
1	B	24	THR	2.4
1	B	28	ILE	2.4
1	C	52	ASP	2.4
1	D	50	ASN	2.3
1	A	48	SER	2.2
1	C	133	ALA	2.2
1	C	53	THR	2.2
1	C	48	SER	2.1
1	B	222	VAL	2.1
1	A	17	ASP	2.1
1	C	16	SER	2.0
1	A	13	ASP	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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
2	GOL	D	301	6/6	0.67	0.16	89,100,103,104	0
2	GOL	C	301	6/6	0.70	0.18	73,89,93,101	0
4	SO4	D	303	5/5	0.72	0.09	142,155,186,193	0
4	SO4	A	308	5/5	0.73	0.10	102,119,149,182	0
3	CL	A	304	1/1	0.74	0.26	123,123,123,123	0
4	SO4	A	305	5/5	0.75	0.09	113,127,148,165	0
2	GOL	C	303	6/6	0.79	0.21	83,95,104,110	0
2	GOL	C	302	6/6	0.79	0.17	88,95,96,103	0
2	GOL	A	303	6/6	0.80	0.14	76,93,98,103	0
2	GOL	A	301	6/6	0.82	0.12	82,91,93,97	0
4	SO4	C	306	5/5	0.86	0.13	69,75,109,126	0
4	SO4	A	307	5/5	0.88	0.14	80,83,104,129	5
4	SO4	A	306	5/5	0.89	0.17	75,75,87,93	5
2	GOL	A	302	6/6	0.90	0.10	61,67,78,82	0
4	SO4	B	302	5/5	0.91	0.10	73,83,100,104	0
4	SO4	B	301	5/5	0.92	0.11	53,66,71,75	0
3	CL	C	304	1/1	0.93	0.07	68,68,68,68	0
4	SO4	C	307	5/5	0.94	0.09	67,68,73,75	0
3	CL	C	305	1/1	0.94	0.06	76,76,76,76	0
3	CL	D	302	1/1	0.95	0.12	62,62,62,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.