



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:52 PM GMT

PDB ID : 4Q7A
Title : Crystal Structure of N-acetyl-ornithine/N-acetyl-lysine Deacetylase from *Sphaerobacter thermophilus*
Authors : Kim, Y.; Tesar, C.; Clancy, S.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2014-04-24
Resolution : 2.05 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

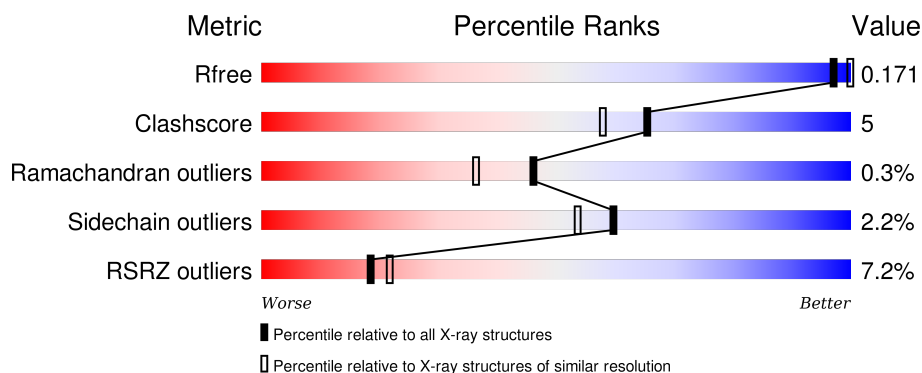
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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}	91344	1192 (2.04-2.04)
Clashscore	102246	1269 (2.04-2.04)
Ramachandran outliers	100387	1258 (2.04-2.04)
Sidechain outliers	100360	1258 (2.04-2.04)
RSRZ outliers	91569	1194 (2.04-2.04)

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. 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	370	<div> <div>5%</div> <div>88%</div> <div>10%</div> <div>•</div> </div>
1	B	370	<div> <div>4%</div> <div>88%</div> <div>10%</div> <div>•</div> </div>
1	C	370	<div> <div>10%</div> <div>79%</div> <div>16%</div> <div>• 5%</div> </div>
1	D	370	<div> <div>8%</div> <div>78%</div> <div>14%</div> <div>• 8%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	C	401	-	-	-	X
3	GOL	D	401	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11552 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.

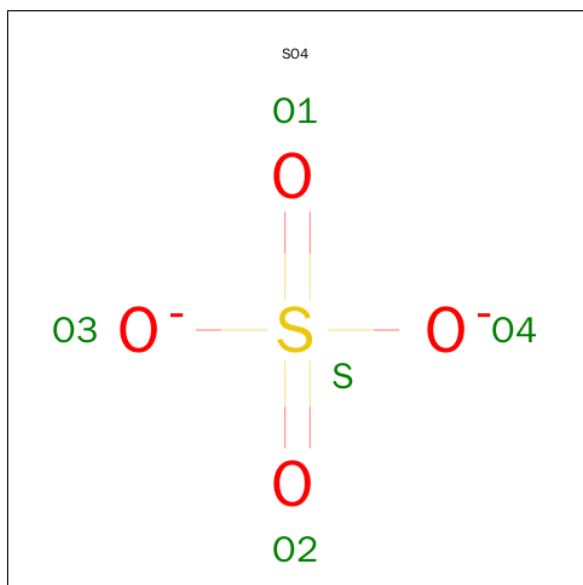
- Molecule 1 is a protein called N-acetyl-ornithine/N-acetyl-lysine deacetylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	361	Total	C	N	O	S	Se	0	0	0
			2649	1664	469	505	3	8			
1	B	361	Total	C	N	O	S	Se	0	2	0
			2670	1678	471	510	3	8			
1	C	352	Total	C	N	O	S	Se	0	1	0
			2591	1628	461	492	3	7			
1	D	341	Total	C	N	O	S	Se	0	2	0
			2529	1589	451	479	3	7			

There are 12 discrepancies between the modelled and reference sequences:

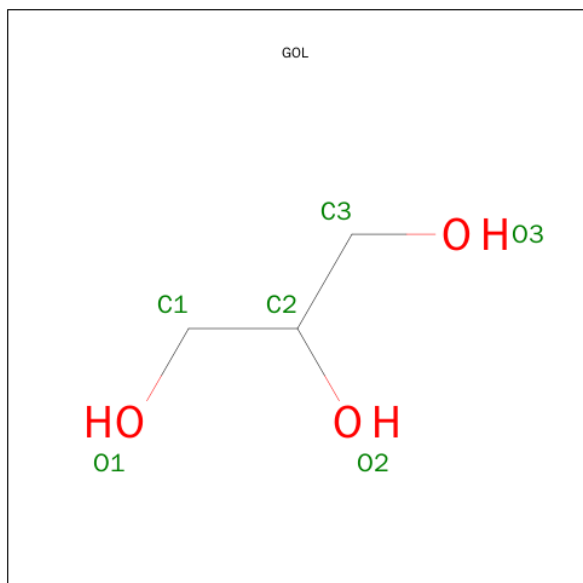
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP D1C6B5
A	-1	ASN	-	EXPRESSION TAG	UNP D1C6B5
A	0	ALA	-	EXPRESSION TAG	UNP D1C6B5
B	-2	SER	-	EXPRESSION TAG	UNP D1C6B5
B	-1	ASN	-	EXPRESSION TAG	UNP D1C6B5
B	0	ALA	-	EXPRESSION TAG	UNP D1C6B5
C	-2	SER	-	EXPRESSION TAG	UNP D1C6B5
C	-1	ASN	-	EXPRESSION TAG	UNP D1C6B5
C	0	ALA	-	EXPRESSION TAG	UNP D1C6B5
D	-2	SER	-	EXPRESSION TAG	UNP D1C6B5
D	-1	ASN	-	EXPRESSION TAG	UNP D1C6B5
D	0	ALA	-	EXPRESSION TAG	UNP D1C6B5

- Molecule 2 is SULFATE ION (three-letter code: 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		
2	B	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	1	Total	Cl	0	0
			1	1		

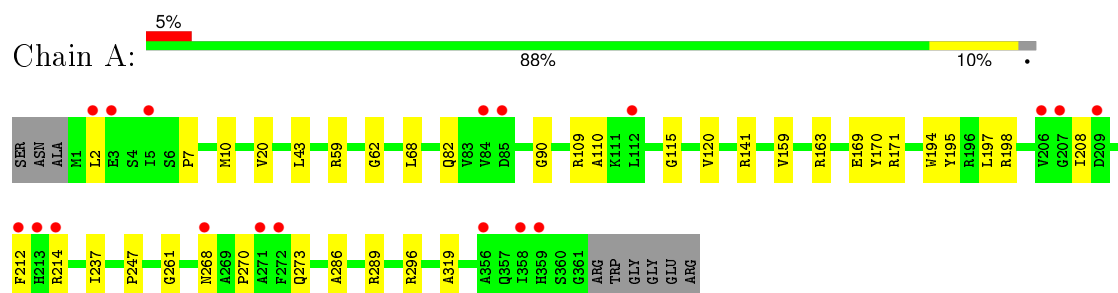
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	273	Total	O	0	0
			273	273		
5	B	269	Total	O	0	0
			269	269		
5	C	242	Total	O	0	0
			242	242		
5	D	284	Total	O	0	0
			284	284		

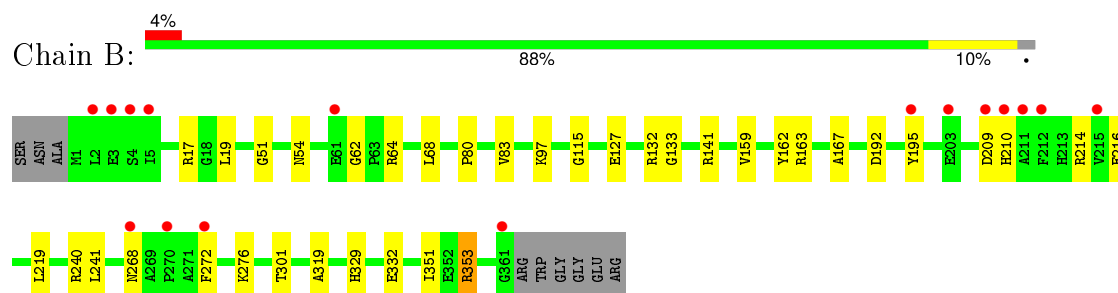
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

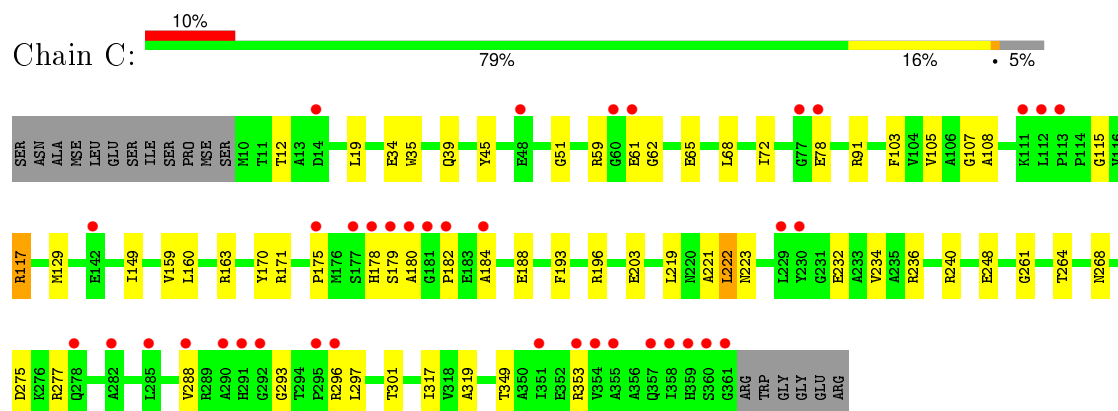
- Molecule 1: N-acetyl-ornithine/N-acetyl-lysine deacetylase



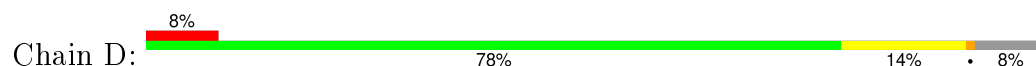
- Molecule 1: N-acetyl-ornithine/N-acetyl-lysine deacetylase

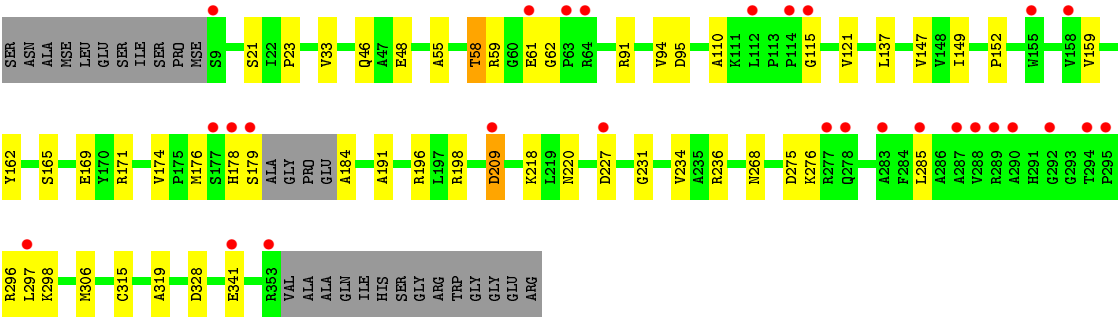


- Molecule 1: N-acetyl-ornithine/N-acetyl-lysine deacetylase



- Molecule 1: N-acetyl-ornithine/N-acetyl-lysine deacetylase





4 Data and refinement statistics

Property	Value	Source
Space group	P 42	Depositor
Cell constants a, b, c, α , β , γ	119.54Å 119.54Å 119.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.78 – 2.05 48.78 – 2.05	Depositor EDS
% Data completeness (in resolution range)	97.0 (48.78-2.05) 97.0 (48.78-2.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.01 (at 2.05Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1678)	Depositor
R, R_{free}	0.154 , 0.171 0.154 , 0.171	Depositor DCC
R_{free} test set	5104 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	24.2	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.6	EDS
Estimated twinning fraction	0.500 for h,-k,-l 0.008 for -h,-l,-k 0.005 for -h,l,k 0.007 for l,-k,h 0.011 for -l,-k,-h 0.487 for h,-k,-l	Xtriage
Reported twinning fraction	0.500 for h,-k,-l	Depositor
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 102060 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11552	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 3.42% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: GOL, SO4, CL

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	0/2701	0.44	0/3681
1	B	0.23	0/2723	0.43	0/3711
1	C	0.27	0/2644	0.45	0/3607
1	D	0.23	0/2578	0.44	0/3512
All	All	0.24	0/10646	0.44	0/14511

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	2649	0	2638	19	0
1	B	2670	0	2651	23	0
1	C	2591	0	2575	36	0
1	D	2529	0	2517	32	0
2	A	5	0	0	0	0
2	B	10	0	0	1	0
2	D	5	0	0	0	0
3	A	6	0	8	0	0
3	C	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	12	0	16	0	0
4	D	1	0	0	0	0
5	A	273	0	0	4	0
5	B	269	0	0	4	0
5	C	242	0	0	6	0
5	D	284	0	0	5	0
All	All	11552	0	10413	105	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 5.

All (105) 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:247:PRO:HG3	1:A:270:PRO:HB3	1.64	0.78
1:C:163:ARG:HH11	1:C:301:THR:HG23	1.48	0.77
1:C:159:VAL:HB	1:C:319:ALA:HB3	1.72	0.72
1:A:10:MSE:O	1:A:109:ARG:NH2	2.24	0.71
1:D:169:GLU:OE2	1:D:236:ARG:NH1	2.25	0.70
1:B:332:GLU:HB2	1:D:48:GLU:HG2	1.73	0.69
1:C:51:GLY:HA2	1:D:179:SER:HB3	1.74	0.68
1:B:163:ARG:HH22	1:B:216:GLU:HG2	1.58	0.68
1:B:219:LEU:HB3	1:C:188:GLU:HG3	1.74	0.67
1:A:159:VAL:HB	1:A:319:ALA:HB3	1.75	0.67
1:C:219:LEU:HD21	1:C:222:LEU:HD12	1.79	0.65
1:C:171:ARG:HG2	1:C:234:VAL:HG22	1.79	0.65
1:B:353:ARG:NH2	5:B:713:HOH:O	2.28	0.65
1:B:159:VAL:HB	1:B:319:ALA:HB3	1.78	0.64
1:D:285:LEU:HD21	1:D:297:LEU:HD21	1.81	0.63
1:D:178:HIS:O	1:D:179:SER:CB	2.47	0.62
1:D:178:HIS:O	1:D:179:SER:HB2	1.99	0.62
1:D:59:ARG:NH2	1:D:110:ALA:O	2.24	0.62
1:A:59:ARG:NH2	5:A:587:HOH:O	2.33	0.61
1:A:169:GLU:OE1	1:A:171:ARG:NE	2.29	0.60
1:D:46:GLN:HG3	1:D:58:THR:HG23	1.83	0.60
1:A:208:ILE:O	1:A:214:ARG:NH1	2.35	0.60
1:C:160:LEU:HD21	1:C:297:LEU:HD22	1.84	0.60
1:B:19:LEU:HG	1:B:97:LYS:HD3	1.83	0.59
1:A:273:GLN:NE2	5:A:613:HOH:O	2.35	0.59
1:D:174:VAL:O	1:D:231:GLY:N	2.28	0.59
1:C:59:ARG:NH2	1:C:107:GLY:O	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:196:ARG:NH1	5:D:620:HOH:O	2.34	0.59
1:D:276:LYS:NZ	5:D:679:HOH:O	2.35	0.58
1:B:127:GLU:OE1	1:B:329:HIS:NE2	2.35	0.57
1:B:163:ARG:NH2	1:B:214:ARG:O	2.37	0.57
1:C:171:ARG:NH1	1:C:232:GLU:OE1	2.37	0.57
1:C:240:ARG:NH1	5:C:615:HOH:O	2.38	0.56
1:D:91:ARG:NH2	5:D:517:HOH:O	2.34	0.56
1:D:198:ARG:NH1	5:D:616:HOH:O	2.38	0.56
1:C:51:GLY:CA	1:D:179:SER:HB3	2.35	0.56
1:D:159:VAL:HB	1:D:319:ALA:HB3	1.87	0.55
1:A:20:VAL:HG11	1:A:90:GLY:HA3	1.89	0.54
1:A:7:PRO:HA	1:A:43:LEU:HD22	1.89	0.54
1:B:167:ALA:HB2	1:B:268:ASN:HD21	1.73	0.53
1:A:68:LEU:HD12	1:A:120:VAL:HG22	1.91	0.53
1:D:62:GLY:HA3	1:D:115:GLY:HA2	1.92	0.51
1:C:61:GLU:HA	1:C:117:ARG:HD3	1.91	0.51
1:A:286:ALA:HA	1:A:289:ARG:HE	1.76	0.51
1:B:240:ARG:HD3	1:B:301:THR:HG23	1.93	0.50
1:B:17:ARG:NH2	1:B:83:VAL:HG21	2.25	0.50
1:C:349:THR:O	1:C:353:ARG:HG2	2.11	0.50
1:C:68:LEU:HB3	1:C:103:PHE:CD2	2.47	0.50
1:A:171:ARG:NH1	5:A:611:HOH:O	2.44	0.50
1:B:162:TYR:CZ	1:B:272:PHE:HB3	2.47	0.49
1:C:240:ARG:NH2	5:C:637:HOH:O	2.34	0.49
1:A:163:ARG:HD2	1:A:212:PHE:HA	1.95	0.49
1:D:171:ARG:HG2	1:D:234:VAL:HG12	1.95	0.48
1:D:94:VAL:HG12	1:D:152:PRO:HG3	1.96	0.47
1:C:65:GLU:HG3	1:C:117:ARG:HB3	1.95	0.47
1:C:184:ALA:HB1	1:C:188:GLU:OE1	2.14	0.47
1:C:178:HIS:O	1:C:180:ALA:N	2.48	0.47
1:D:165:SER:HB3	1:D:268:ASN:HD22	1.80	0.47
1:B:192:ASP:HA	1:B:195[A]:TYR:CD1	2.50	0.46
1:B:54:ASN:HD21	1:B:133:GLY:HA2	1.81	0.46
1:A:195:TYR:OH	1:D:191:ALA:O	2.26	0.46
1:D:209:ASP:N	1:D:209:ASP:OD1	2.47	0.46
1:C:163:ARG:NH1	1:C:301:THR:HG23	2.26	0.46
1:C:240:ARG:NH1	5:C:582:HOH:O	2.49	0.45
1:B:62:GLY:HA3	1:B:115:GLY:HA2	1.97	0.45
1:C:248:GLU:OE2	5:C:578:HOH:O	2.21	0.45
1:B:68:LEU:HD11	1:B:351:ILE:HD11	1.98	0.45
1:C:35:TRP:O	1:C:39:GLN:HG2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:GLY:N	2:B:402:SO4:O1	2.44	0.45
1:C:170:TYR:CZ	1:C:261:GLY:HA3	2.52	0.45
1:D:296:ARG:HG3	1:D:298:LYS:HE2	2.00	0.45
1:D:147:VAL:HG23	1:D:315:CYS:SG	2.57	0.45
1:B:17:ARG:HH21	1:B:83:VAL:HG21	1.83	0.44
1:D:149:ILE:HG12	1:D:306:MSE:HB2	2.00	0.43
1:C:62:GLY:HA3	1:C:115:GLY:HA2	2.00	0.43
1:C:193:PHE:HA	1:C:196:ARG:NH1	2.33	0.43
1:C:129[A]:MSE:SE	1:C:268:ASN:HB2	2.69	0.43
1:D:169:GLU:OE1	1:D:171:ARG:NH2	2.47	0.43
1:C:221:ALA:HB3	1:C:236:ARG:HE	1.84	0.43
1:C:91:ARG:O	5:C:589:HOH:O	2.21	0.43
1:B:141:ARG:O	5:B:599:HOH:O	2.22	0.42
1:C:203:GLU:OE1	5:C:547:HOH:O	2.22	0.42
1:D:21:SER:O	1:D:23:PRO:HD3	2.19	0.42
1:D:33:VAL:HG13	1:D:55:ALA:HB2	2.01	0.42
1:B:163:ARG:HB3	1:B:241:LEU:O	2.19	0.42
1:C:288:VAL:HG12	1:C:293:GLY:HA3	2.01	0.42
1:D:121:VAL:HG11	1:D:137:LEU:HD11	2.02	0.42
1:C:296:ARG:HB2	1:C:296:ARG:HE	1.67	0.42
1:A:170:TYR:CZ	1:A:261:GLY:HA3	2.54	0.42
1:B:132:ARG:NH1	5:B:652:HOH:O	2.53	0.42
1:B:17:ARG:NH1	5:B:703:HOH:O	2.53	0.41
1:C:45:TYR:CE1	1:C:108:ALA:HB2	2.54	0.41
1:C:223:ASN:HD22	1:C:236:ARG:NH2	2.17	0.41
1:B:64:ARG:HA	1:B:64:ARG:HD3	1.70	0.41
1:D:218:LYS:HE2	1:D:220:ASN:HD21	1.86	0.41
1:A:62:GLY:HA3	1:A:115:GLY:HA2	2.02	0.41
1:A:194:TRP:CZ2	1:A:198:ARG:HD3	2.56	0.41
1:A:110:ALA:O	5:A:587:HOH:O	2.22	0.41
1:C:19:LEU:HD11	1:C:72:ILE:HD11	2.02	0.41
1:D:159:VAL:HG11	1:D:162:TYR:CZ	2.55	0.41
1:A:197:LEU:HD23	1:A:237:ILE:HG21	2.03	0.41
1:C:12:THR:HA	1:C:105:VAL:HG21	2.02	0.41
1:D:95:ASP:HB2	1:D:328:ASP:OD1	2.21	0.40
1:C:149:ILE:HD11	1:C:317:ILE:HD11	2.04	0.40
1:D:184:ALA:N	5:D:710:HOH:O	2.53	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	359/370 (97%)	345 (96%)	14 (4%)	0	100	100
1	B	361/370 (98%)	349 (97%)	11 (3%)	1 (0%)	46	36
1	C	351/370 (95%)	334 (95%)	14 (4%)	3 (1%)	21	10
1	D	339/370 (92%)	326 (96%)	13 (4%)	0	100	100
All	All	1410/1480 (95%)	1354 (96%)	52 (4%)	4 (0%)	46	36

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	179	SER
1	C	182	PRO
1	C	175	PRO
1	B	80	PRO

5.3.2 Protein sidechains [i](#)

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	271/269 (101%)	266 (98%)	5 (2%)	66	62
1	B	273/269 (102%)	269 (98%)	4 (2%)	72	70
1	C	263/269 (98%)	256 (97%)	7 (3%)	52	45
1	D	258/269 (96%)	251 (97%)	7 (3%)	52	45
All	All	1065/1076 (99%)	1042 (98%)	23 (2%)	60	53

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

Mol	Chain	Res	Type
1	A	2	LEU
1	A	82	GLN
1	A	141	ARG
1	A	268	ASN
1	A	296	ARG
1	B	209	ASP
1	B	210	HIS
1	B	276	LYS
1	B	353	ARG
1	C	34	GLU
1	C	78	GLU
1	C	117	ARG
1	C	222	LEU
1	C	264	THR
1	C	275	ASP
1	C	277	ARG
1	D	58	THR
1	D	61	GLU
1	D	176	MSE
1	D	209	ASP
1	D	227	ASP
1	D	275	ASP
1	D	341	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	223	ASN
1	B	39	GLN
1	B	223	ASN
1	B	268	ASN
1	C	82	GLN
1	C	213	HIS
1	C	329	HIS
1	D	46	GLN
1	D	82	GLN
1	D	220	ASN
1	D	268	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 9 ligands modelled in this entry, 1 is 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	$\# Z > 2$
2	SO4	A	401	-	4,4,4	0.32	0	6,6,6	0.10	0
3	GOL	A	402	-	5,5,5	0.32	0	5,5,5	0.24	0
2	SO4	B	401	-	4,4,4	0.30	0	6,6,6	0.11	0
2	SO4	B	402	-	4,4,4	0.35	0	6,6,6	0.12	0
3	GOL	C	401	-	5,5,5	0.32	0	5,5,5	0.40	0
3	GOL	D	401	-	5,5,5	0.35	0	5,5,5	0.25	0
2	SO4	D	402	-	4,4,4	0.32	0	6,6,6	0.09	0
3	GOL	D	403	-	5,5,5	0.33	0	5,5,5	0.26	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	SO4	A	401	-	-	0/0/0/0	0/0/0/0
3	GOL	A	402	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	B	401	-	-	0/0/0/0	0/0/0/0
2	SO4	B	402	-	-	0/0/0/0	0/0/0/0
3	GOL	C	401	-	-	0/4/4/4	0/0/0/0
3	GOL	D	401	-	-	0/4/4/4	0/0/0/0
2	SO4	D	402	-	-	0/0/0/0	0/0/0/0
3	GOL	D	403	-	-	0/4/4/4	0/0/0/0

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
2	B	402	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	353/370 (95%)	0.57	18 (5%) 32 36	11, 25, 42, 66	0
1	B	353/370 (95%)	0.57	16 (4%) 37 42	12, 24, 42, 66	0
1	C	346/370 (93%)	0.81	38 (10%) 7 8	14, 26, 48, 91	0
1	D	335/370 (90%)	0.73	28 (8%) 14 15	13, 26, 45, 73	1 (0%)
All	All	1387/1480 (93%)	0.67	100 (7%) 18 22	11, 25, 45, 91	1 (0%)

All (100) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	182	PRO	10.7
1	C	177	SER	9.4
1	C	178	HIS	7.8
1	C	179	SER	7.0
1	C	181	GLY	5.9
1	C	180	ALA	5.3
1	D	178	HIS	5.3
1	C	282	ALA	5.1
1	C	358	ILE	4.9
1	C	112	LEU	4.8
1	D	179	SER	4.7
1	B	2	LEU	4.6
1	B	4	SER	4.4
1	D	290	ALA	4.3
1	C	360	SER	4.3
1	D	9	SER	4.2
1	C	357	GLN	4.1
1	D	353	ARG	3.9
1	B	272	PHE	3.8
1	A	207	GLY	3.7
1	C	290	ALA	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	2	LEU	3.7
1	A	268	ASN	3.6
1	C	184	ALA	3.6
1	D	115	GLY	3.6
1	C	288	VAL	3.5
1	D	292	GLY	3.5
1	D	277	ARG	3.5
1	B	361	GLY	3.5
1	B	3	GLU	3.4
1	D	288	VAL	3.3
1	A	112	LEU	3.3
1	A	3	GLU	3.3
1	D	227	ASP	3.3
1	D	287	ALA	3.1
1	C	230	TYR	3.1
1	C	359	HIS	3.1
1	D	294	THR	3.1
1	A	5	ILE	3.1
1	C	354	VAL	3.0
1	D	158	VAL	3.0
1	B	209	ASP	3.0
1	C	285	LEU	2.9
1	C	361	GLY	2.9
1	A	272	PHE	2.8
1	D	177	SER	2.8
1	C	142	GLU	2.8
1	C	351	ILE	2.8
1	D	341	GLU	2.7
1	C	229	LEU	2.7
1	C	292	GLY	2.7
1	D	297	LEU	2.7
1	A	359	HIS	2.7
1	C	14	ASP	2.6
1	D	155	TRP	2.6
1	B	195[A]	TYR	2.6
1	A	85	ASP	2.6
1	C	278	GLN	2.6
1	B	5	ILE	2.6
1	A	271	ALA	2.5
1	C	355	ALA	2.5
1	B	215	VAL	2.4
1	C	78	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	111	LYS	2.4
1	B	203	GLU	2.4
1	A	206	VAL	2.4
1	D	283	ALA	2.4
1	C	353	ARG	2.4
1	D	61	GLU	2.4
1	A	214	ARG	2.4
1	B	268	ASN	2.4
1	D	278	GLN	2.3
1	A	212	PHE	2.3
1	C	295	PRO	2.3
1	D	289	ARG	2.3
1	D	112	LEU	2.3
1	A	213	HIS	2.3
1	D	285	LEU	2.2
1	B	212	PHE	2.2
1	B	61	GLU	2.2
1	C	61	GLU	2.2
1	C	48	GLU	2.2
1	A	84	VAL	2.2
1	D	64	ARG	2.1
1	B	210	HIS	2.1
1	A	209	ASP	2.1
1	B	211	ALA	2.1
1	B	270	PRO	2.1
1	C	77	GLY	2.1
1	C	175	PRO	2.1
1	D	63	PRO	2.1
1	A	356	ALA	2.1
1	C	113	PRO	2.0
1	D	114	PRO	2.0
1	C	296	ARG	2.0
1	D	295	PRO	2.0
1	A	358	ILE	2.0
1	D	209	ASP	2.0
1	C	291	HIS	2.0
1	C	60	GLY	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 [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	GOL	D	401	6/6	0.83	0.33	9.50	28,32,40,49	0
3	GOL	C	401	6/6	0.67	0.30	7.91	39,50,55,56	0
2	SO4	D	402	5/5	0.93	0.22	1.51	44,47,53,60	0
2	SO4	A	401	5/5	0.92	0.16	0.93	33,41,50,56	0
2	SO4	B	401	5/5	0.88	0.20	-0.05	42,45,46,48	0
3	GOL	D	403	6/6	0.72	0.42	-	40,43,43,45	0
2	SO4	B	402	5/5	0.94	0.20	-	37,39,45,48	0
3	GOL	A	402	6/6	0.69	0.26	-	48,51,55,55	0
4	CL	D	404	1/1	0.81	0.12	-	46,46,46,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.