Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 03:53 pm BST

PDB ID	:	2YFQ
Title	:	Crystal structure of Glutamate dehydrogenase from Peptoniphilus asaccharolyticus
Authors	:	Oliveira, T.; Panjikar, S.; Carrigan, J.B.; Sharkey, M.A.; Hamza, M.; Engel, P.C.; Khan, A.R.
Deposited on	:	2011-04-07
Resolution	:	2.94 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:

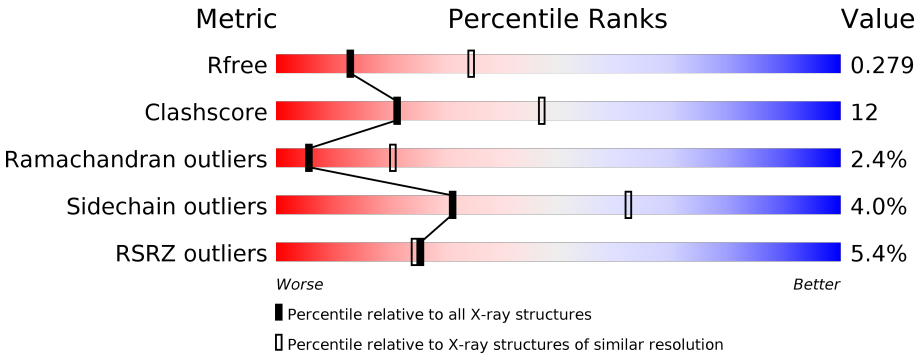
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.94 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	2969 (2.98-2.90)
Clashscore	141614	3218 (2.98-2.90)
Ramachandran outliers	138981	3122 (2.98-2.90)
Sidechain outliers	138945	3124 (2.98-2.90)
RSRZ outliers	127900	2902 (2.98-2.90)

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 <=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	421	5% 68% 23% • 6%
1	B	421	5% 68% 22% • 6%

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
2	SO4	B	1423	-	-	X	-

2 Entry composition

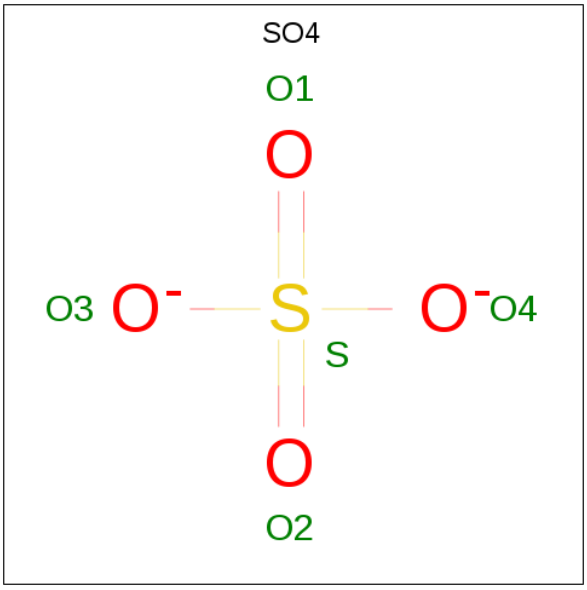
There are 3 unique types of molecules in this entry. The entry contains 6108 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 NAD-SPECIFIC GLUTAMATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	395	Total	C	N	O	S	0	0	0
			3033	1925	516	577	15			
1	B	397	Total	C	N	O	S	0	0	0
			3049	1935	518	581	15			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

- Molecule 3 is water.

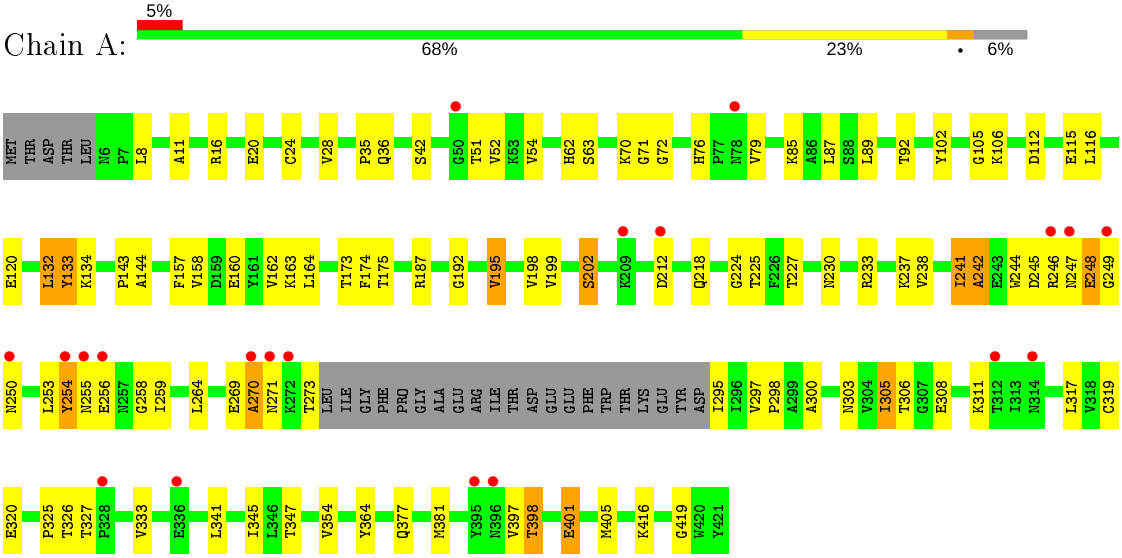
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	4	Total O 4 4	0	0
3	B	2	Total O 2 2	0	0

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 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.

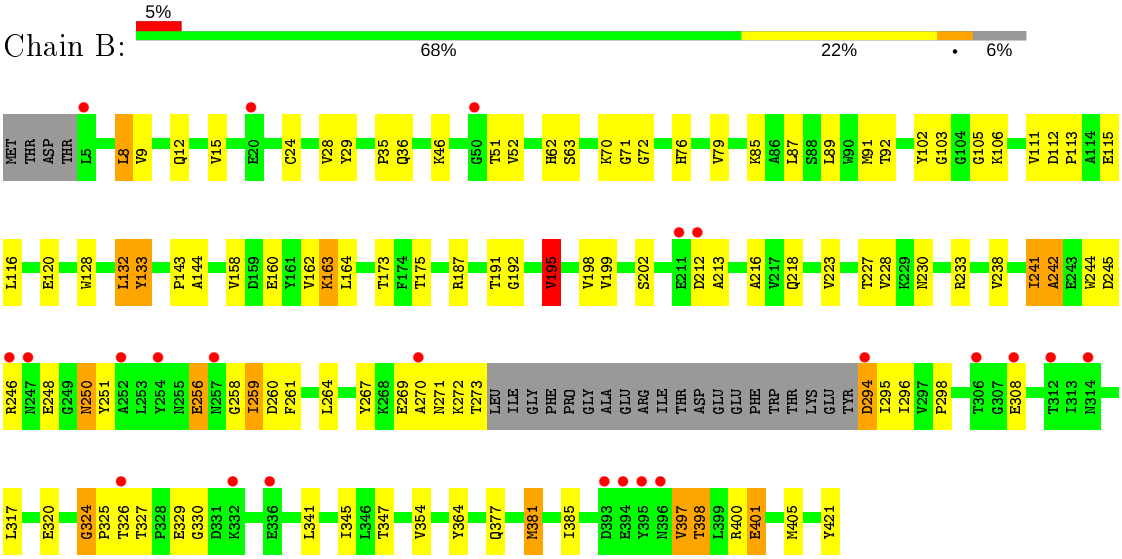
- Molecule 1: NAD-SPECIFIC GLUTAMATE DEHYDROGENASE

Chain A:



- Molecule 1: NAD-SPECIFIC GLUTAMATE DEHYDROGENASE

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α, β, γ	153.31Å 153.31Å 318.99Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.94 19.74 – 2.94	Depositor EDS
% Data completeness (in resolution range)	97.8 (20.00-2.94) 91.2 (19.74-2.94)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.59 (at 2.93Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.242 , 0.287 0.241 , 0.279	Depositor DCC
R_{free} test set	935 reflections (3.33%)	wwPDB-VP
Wilson B-factor (Å^2)	63.6	Xtriage
Anisotropy	0.002	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 75.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	6108	wwPDB-VP
Average B, all atoms (Å^2)	79.0	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 70.89 % 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 2.8888e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.63	5/3092 (0.2%)	0.70	1/4183 (0.0%)
1	B	0.63	6/3108 (0.2%)	0.69	1/4205 (0.0%)
All	All	0.63	11/6200 (0.2%)	0.70	2/8388 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	2
All	All	0	3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	133	TYR	CE2-CZ	-7.69	1.28	1.38
1	A	133	TYR	CE1-CZ	-7.09	1.29	1.38
1	B	195	VAL	CB-CG1	-6.83	1.38	1.52
1	A	195	VAL	CB-CG2	-6.78	1.38	1.52
1	A	133	TYR	CE2-CZ	-6.47	1.30	1.38
1	B	133	TYR	CG-CD1	-6.01	1.31	1.39
1	B	195	VAL	CB-CG2	-5.84	1.40	1.52
1	B	133	TYR	CE1-CZ	-5.83	1.30	1.38
1	A	133	TYR	CG-CD1	-5.72	1.31	1.39
1	B	133	TYR	CG-CD2	-5.71	1.31	1.39
1	A	133	TYR	CG-CD2	-5.63	1.31	1.39

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	246	ARG	N-CA-C	5.11	124.80	111.00
1	B	246	ARG	N-CA-C	5.11	124.80	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	255	ASN	Peptide
1	B	294	ASP	Peptide
1	B	324	GLY	Peptide

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	3033	0	2992	68	0
1	B	3049	0	3007	80	0
2	A	10	0	0	1	0
2	B	10	0	0	3	0
3	A	4	0	0	0	0
3	B	2	0	0	0	0
All	All	6108	0	5999	145	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 12.

All (145) 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:241:ILE:HG12	1:A:259:ILE:HD12	1.41	1.03
1:B:87:LEU:HB3	1:B:106:LYS:HD3	1.48	0.95
1:B:187:ARG:NH1	2:B:1422:SO4:O1	2.00	0.94
1:B:294:ASP:HB3	1:B:295:ILE:HG23	1.52	0.91
1:A:87:LEU:HB3	1:A:106:LYS:HD3	1.63	0.80
1:B:8:LEU:HD11	1:B:85:LYS:HB3	1.68	0.75
1:A:187:ARG:NH1	2:A:1423:SO4:O2	2.21	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:VAL:CG1	1:B:227:THR:HG22	2.19	0.71
1:B:62:HIS:HD2	1:B:92:THR:OG1	1.74	0.71
1:B:36:GLN:NE2	1:B:63:SER:HA	2.06	0.70
1:A:241:ILE:CG1	1:A:259:ILE:HD12	2.19	0.70
1:B:195:VAL:HG13	1:B:227:THR:HG22	1.75	0.69
1:B:8:LEU:HD11	1:B:85:LYS:CB	2.23	0.67
1:B:327:THR:O	1:B:330:GLY:N	2.27	0.66
1:A:354:VAL:HG23	1:A:377:GLN:OE1	1.96	0.66
1:A:230:ASN:HA	1:A:233:ARG:HB2	1.78	0.65
1:B:267:TYR:O	1:B:271:ASN:HB2	1.96	0.65
1:B:223:VAL:O	1:B:227:THR:HG23	1.98	0.64
1:B:112:ASP:O	1:B:115:GLU:HG2	1.97	0.63
1:B:250:ASN:H	1:B:250:ASN:HD22	1.47	0.63
1:A:295:ILE:HB	1:A:317:LEU:O	2.00	0.62
1:A:192:GLY:HA2	1:A:227:THR:HG22	1.81	0.62
1:B:87:LEU:CB	1:B:106:LYS:HD3	2.26	0.61
1:B:230:ASN:HA	1:B:233:ARG:HB2	1.82	0.61
1:B:87:LEU:HB3	1:B:106:LYS:CD	2.26	0.61
1:A:87:LEU:HB3	1:A:106:LYS:CD	2.30	0.60
1:A:198:VAL:O	1:A:202:SER:OG	2.19	0.60
1:B:327:THR:HB	1:B:329:GLU:HG3	1.84	0.59
1:A:36:GLN:NE2	1:A:63:SER:HA	2.18	0.59
1:A:133:TYR:CD2	1:A:164:LEU:HD13	2.37	0.59
1:A:218:GLN:HE21	1:A:300:ALA:HB3	1.69	0.58
1:A:62:HIS:HD2	1:A:92:THR:OG1	1.85	0.58
1:A:305:ILE:HG21	1:A:326:THR:HG22	1.84	0.58
1:B:354:VAL:HG23	1:B:377:GLN:OE1	2.04	0.57
1:A:238:VAL:O	1:A:258:GLY:HA2	2.04	0.57
1:A:377:GLN:O	1:A:381:MET:HG2	2.04	0.57
1:A:112:ASP:O	1:A:115:GLU:HG2	2.04	0.57
1:A:16:ARG:O	1:A:20:GLU:HB2	2.05	0.56
1:A:87:LEU:CB	1:A:106:LYS:HD3	2.33	0.56
1:A:253:LEU:O	1:A:254:TYR:HB2	2.05	0.56
1:A:269:GLU:O	1:A:270:ALA:HB3	2.07	0.55
1:A:160:GLU:OE1	1:A:160:GLU:HA	2.07	0.54
1:B:71:GLY:HA3	1:B:105:GLY:O	2.08	0.53
1:B:398:THR:HB	1:B:401:GLU:HG2	1.89	0.53
1:A:241:ILE:O	1:A:242:ALA:CB	2.57	0.53
1:A:51:THR:HG22	1:A:52:VAL:N	2.24	0.53
1:B:76:HIS:O	1:B:79:VAL:HG12	2.08	0.53
1:A:158:VAL:O	1:A:162:VAL:HG23	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:HIS:O	1:A:79:VAL:HG12	2.09	0.53
1:A:241:ILE:HG12	1:A:259:ILE:CD1	2.28	0.52
1:B:198:VAL:O	1:B:202:SER:OG	2.23	0.52
1:B:195:VAL:HG11	1:B:227:THR:CG2	2.39	0.52
1:B:241:ILE:O	1:B:242:ALA:CB	2.57	0.52
1:B:326:THR:OG1	1:B:400:ARG:NH2	2.42	0.52
1:A:71:GLY:HA3	1:A:105:GLY:O	2.11	0.51
1:B:195:VAL:CG1	1:B:227:THR:CG2	2.87	0.51
1:A:259:ILE:HG21	1:A:264:LEU:HD22	1.92	0.51
1:B:216:ALA:HB3	1:B:296:ILE:HG12	1.92	0.51
1:A:116:LEU:HD22	1:A:120:GLU:HB3	1.93	0.51
1:B:160:GLU:OE1	1:B:160:GLU:HA	2.10	0.51
1:A:398:THR:HB	1:A:401:GLU:HG2	1.94	0.50
1:B:15:VAL:HG12	1:B:29:TYR:HD1	1.76	0.50
1:B:143:PRO:HD2	1:B:173:THR:O	2.12	0.50
1:B:116:LEU:HD22	1:B:120:GLU:HB3	1.94	0.50
1:A:305:ILE:CG2	1:A:326:THR:HG22	2.41	0.50
1:B:213:ALA:HB1	1:B:294:ASP:HB2	1.94	0.50
1:A:397:VAL:HG22	1:A:398:THR:H	1.77	0.49
1:B:51:THR:HG22	1:B:52:VAL:N	2.27	0.49
1:A:244:TRP:CE2	1:A:245:ASP:O	2.65	0.49
1:B:401:GLU:O	1:B:405:MET:SD	2.71	0.49
1:A:70:LYS:HD3	1:A:102:TYR:CE2	2.48	0.48
1:B:213:ALA:HB1	1:B:294:ASP:CG	2.34	0.48
1:B:238:VAL:O	1:B:258:GLY:O	2.31	0.48
1:B:218:GLN:HA	1:B:242:ALA:HB3	1.94	0.48
1:B:271:ASN:C	1:B:273:THR:H	2.16	0.48
1:A:311:LYS:HG2	1:A:333:VAL:HG21	1.95	0.48
1:B:244:TRP:CE2	1:B:245:ASP:O	2.66	0.48
1:B:397:VAL:HG22	1:B:398:THR:H	1.79	0.48
1:A:35:PRO:HG2	1:A:85:LYS:HG2	1.95	0.48
1:B:192:GLY:HA2	1:B:227:THR:HG22	1.96	0.47
1:B:72:GLY:HA2	1:B:144:ALA:O	2.14	0.47
1:A:269:GLU:O	1:A:270:ALA:CB	2.62	0.47
1:B:320:GLU:O	1:B:347:THR:HB	2.14	0.47
1:B:191:THR:O	1:B:195:VAL:HG12	2.15	0.47
1:B:62:HIS:CD2	1:B:92:THR:OG1	2.61	0.47
1:A:297:VAL:HG22	1:A:319:CYS:HB2	1.96	0.46
1:B:158:VAL:O	1:B:162:VAL:HG23	2.14	0.46
1:A:70:LYS:NZ	1:A:175:THR:OG1	2.37	0.46
1:B:70:LYS:NZ	1:B:175:THR:OG1	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:LYS:HD3	1:B:102:TYR:CE2	2.50	0.46
1:A:419:GLY:O	1:B:46:LYS:NZ	2.39	0.46
1:B:218:GLN:HB3	1:B:298:PRO:HA	1.98	0.46
1:B:327:THR:CB	1:B:329:GLU:HG3	2.44	0.46
1:B:269:GLU:O	1:B:270:ALA:HB3	2.15	0.46
1:A:132:LEU:O	1:A:134:LYS:N	2.49	0.46
1:A:72:GLY:HA2	1:A:144:ALA:O	2.15	0.46
1:A:218:GLN:HB3	1:A:298:PRO:HA	1.98	0.46
1:A:224:GLY:H	1:A:227:THR:HG23	1.80	0.46
1:A:244:TRP:HD1	1:A:250:ASN:H	1.60	0.46
1:A:320:GLU:O	1:A:347:THR:HB	2.15	0.46
1:B:341:LEU:HD21	1:B:400:ARG:HH21	1.81	0.46
1:A:11:ALA:HB1	1:A:89:LEU:HG	1.97	0.45
1:B:91:MET:CE	2:B:1423:SO4:O1	2.65	0.45
1:A:416:LYS:HG3	1:A:416:LYS:O	2.16	0.45
1:B:63:SER:O	1:B:103:GLY:HA3	2.16	0.45
1:B:8:LEU:HD11	1:B:85:LYS:HB2	1.98	0.45
1:A:132:LEU:O	1:A:133:TYR:C	2.55	0.45
1:A:195:VAL:O	1:A:199:VAL:HG23	2.18	0.44
1:B:91:MET:HE1	2:B:1423:SO4:O1	2.18	0.44
1:A:143:PRO:HD2	1:A:173:THR:O	2.17	0.44
1:B:377:GLN:O	1:B:381:MET:HG2	2.17	0.44
1:A:401:GLU:HG2	1:A:401:GLU:H	1.65	0.44
1:B:9:VAL:HA	1:B:12:GLN:HE21	1.82	0.44
1:B:320:GLU:HG3	1:B:341:LEU:HD11	1.99	0.44
1:A:247:ASN:O	1:A:248:GLU:CB	2.66	0.44
1:A:295:ILE:HD12	1:A:295:ILE:C	2.38	0.43
1:A:42:SER:HB3	1:A:54:VAL:HG11	2.00	0.43
1:B:128:TRP:CE2	1:B:132:LEU:HD11	2.53	0.43
1:B:295:ILE:HD12	1:B:295:ILE:C	2.39	0.43
1:A:52:VAL:HG21	1:B:421:TYR:CB	2.48	0.43
1:B:198:VAL:HG22	1:B:385:ILE:HD11	2.01	0.43
1:B:195:VAL:O	1:B:199:VAL:HG23	2.18	0.43
1:B:24:CYS:HB2	1:B:28:VAL:HG21	2.01	0.43
1:B:242:ALA:HA	1:B:251:TYR:O	2.18	0.43
1:A:271:ASN:C	1:A:273:THR:H	2.21	0.43
1:B:228:VAL:CG2	1:B:261:PHE:HD1	2.31	0.43
1:A:52:VAL:HG21	1:B:421:TYR:HB2	2.00	0.43
1:B:195:VAL:HG11	1:B:227:THR:HG21	2.00	0.43
1:A:163:LYS:HD3	1:A:163:LYS:HA	1.72	0.42
1:B:250:ASN:H	1:B:250:ASN:ND2	2.14	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:306:THR:HA	1:A:327:THR:OG1	2.20	0.42
1:B:295:ILE:HB	1:B:317:LEU:O	2.19	0.42
1:A:247:ASN:O	1:A:248:GLU:HB2	2.18	0.42
1:B:258:GLY:O	1:B:259:ILE:CB	2.68	0.42
1:A:401:GLU:O	1:A:405:MET:SD	2.77	0.42
1:A:24:CYS:HB2	1:A:28:VAL:HG21	2.02	0.41
1:B:163:LYS:HA	1:B:163:LYS:HD3	1.71	0.41
1:B:35:PRO:HA	1:B:62:HIS:HA	2.01	0.41
1:A:320:GLU:HG3	1:A:341:LEU:HD11	2.02	0.41
1:B:324:GLY:O	1:B:326:THR:N	2.52	0.41
1:B:133:TYR:CD1	1:B:164:LEU:HD13	2.55	0.41
1:A:157:PHE:HB3	1:A:174:PHE:CZ	2.56	0.41
1:B:35:PRO:HG2	1:B:85:LYS:HG2	2.02	0.40
1:A:224:GLY:O	1:A:225:THR:C	2.58	0.40
1:B:111:VAL:HG13	1:B:113:PRO:HD3	2.02	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	391/421 (93%)	354 (90%)	26 (7%)	11 (3%)	5 17
1	B	393/421 (93%)	358 (91%)	27 (7%)	8 (2%)	7 25
All	All	784/842 (93%)	712 (91%)	53 (7%)	19 (2%)	6 21

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	242	ALA
1	A	248	GLU
1	A	254	TYR

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Mol	Chain	Res	Type
1	A	256	GLU
1	B	242	ALA
1	A	241	ILE
1	A	270	ALA
1	A	325	PRO
1	B	241	ILE
1	B	325	PRO
1	B	256	GLU
1	B	272	LYS
1	A	305	ILE
1	B	259	ILE
1	A	249	GLY
1	A	303	ASN
1	B	248	GLU
1	B	345	ILE
1	A	345	ILE

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	312/340 (92%)	303 (97%)	9 (3%)	42 73
1	B	314/340 (92%)	298 (95%)	16 (5%)	24 53
All	All	626/680 (92%)	601 (96%)	25 (4%)	31 62

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LEU
1	A	132	LEU
1	A	202	SER
1	A	212	ASP
1	A	237	LYS
1	A	308	GLU
1	A	364	TYR

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Mol	Chain	Res	Type
1	A	398	THR
1	A	401	GLU
1	B	8	LEU
1	B	89	LEU
1	B	132	LEU
1	B	163	LYS
1	B	195	VAL
1	B	212	ASP
1	B	250	ASN
1	B	256	GLU
1	B	260	ASP
1	B	264	LEU
1	B	308	GLU
1	B	364	TYR
1	B	381	MET
1	B	397	VAL
1	B	398	THR
1	B	401	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	36	GLN
1	A	62	HIS
1	A	123	GLN
1	A	218	GLN
1	A	362	ASN
1	B	12	GLN
1	B	36	GLN
1	B	62	HIS
1	B	123	GLN
1	B	250	ASN
1	B	362	ASN

5.3.3 RNA

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 carbohydrates 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	SO4	A	1422	-	4,4,4	0.33	0	6,6,6	0.23	0
2	SO4	B	1422	-	4,4,4	0.31	0	6,6,6	0.37	0
2	SO4	A	1423	-	4,4,4	0.32	0	6,6,6	0.63	0
2	SO4	B	1423	-	4,4,4	0.28	0	6,6,6	0.38	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.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1422	SO4	1	0
2	A	1423	SO4	1	0
2	B	1423	SO4	2	0

5.7 Other polymers

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, 95^{th} 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	395/421 (93%)	-0.13	20 (5%)	28	27	37, 71, 128, 160	0
1	B	397/421 (94%)	-0.00	23 (5%)	23	21	41, 76, 141, 228	0
All	All	792/842 (94%)	-0.07	43 (5%)	25	24	37, 74, 132, 228	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	308	GLU	5.7
1	A	250	ASN	3.8
1	B	254	TYR	3.7
1	A	249	GLY	3.6
1	A	270	ALA	3.5
1	B	270	ALA	3.5
1	A	312	THR	3.5
1	B	326	THR	3.5
1	B	246	ARG	3.4
1	A	271	ASN	3.3
1	B	247	ASN	3.2
1	B	252	ALA	3.2
1	B	336	GLU	3.2
1	A	212	ASP	3.1
1	B	314	ASN	3.1
1	A	314	ASN	3.0
1	B	5	LEU	2.9
1	B	332	LYS	2.9
1	B	211	GLU	2.8
1	A	50	GLY	2.7
1	A	247	ASN	2.7
1	A	395	TYR	2.6
1	B	257	ASN	2.4
1	B	312	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	393	ASP	2.4
1	B	212	ASP	2.4
1	A	255	ASN	2.3
1	A	336	GLU	2.2
1	A	209	LYS	2.2
1	A	328	PRO	2.2
1	B	50	GLY	2.2
1	B	294	ASP	2.2
1	A	246	ARG	2.2
1	A	396	ASN	2.2
1	A	254	TYR	2.2
1	A	256	GLU	2.1
1	B	306	THR	2.1
1	A	272	LYS	2.1
1	B	394	GLU	2.1
1	B	20	GLU	2.1
1	B	395	TYR	2.0
1	B	396	ASN	2.0
1	A	78	ASN	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 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. The B-factors column lists the minimum, median, 95^{th} 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	SO4	A	1423	5/5	0.94	0.19	65,67,68,70	0
2	SO4	B	1422	5/5	0.95	0.21	61,65,66,67	0
2	SO4	A	1422	5/5	0.95	0.19	64,64,66,67	0
2	SO4	B	1423	5/5	0.96	0.17	63,67,68,68	0

6.5 Other polymers (i)

There are no such residues in this entry.