



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 08:48 pm GMT

PDB ID : 1YW6
Title : Crystal Structure of Succinylglutamate Desuccinylase from Escherichia coli, Northeast Structural Genomics Target ET72.
Authors : Forouhar, F.; Yong, W.; Kuzin, A.P.; Ciano, M.; Acton, T.B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2005-02-17
Resolution : 3.10 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

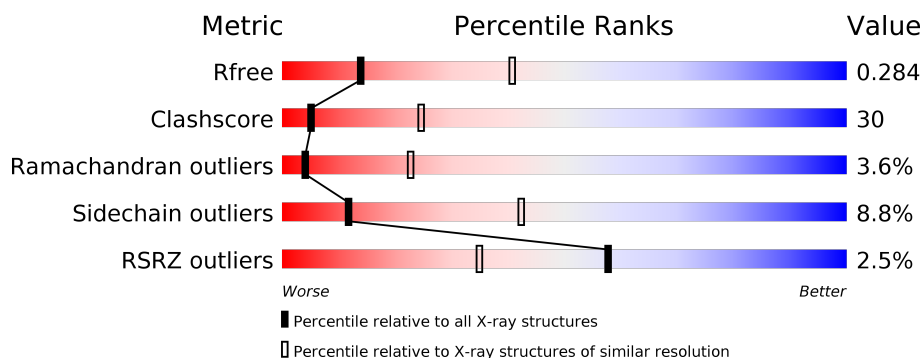
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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}	100719	1001 (3.12-3.08)
Clashscore	112137	1099 (3.12-3.08)
Ramachandran outliers	110173	1057 (3.12-3.08)
Sidechain outliers	110143	1057 (3.12-3.08)
RSRZ outliers	101464	1006 (3.12-3.08)

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	335	<div> <div> <div></div> <div>53%</div> <div>36%</div> <div>5%</div> <div>6%</div> </div> <div>4%</div> </div>
1	B	335	<div> <div> <div></div> <div>39%</div> <div>46%</div> <div>5%</div> <div>10%</div> </div> <div>4%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4969 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 Succinylglutamate desuccinylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	316	Total	C	N	O	S	Se	0	0	0
			2468	1579	435	443	3	8			
1	B	303	Total	C	N	O	S	Se	0	0	0
			2377	1526	419	421	4	7			

There are 42 discrepancies between the modelled and reference sequences:

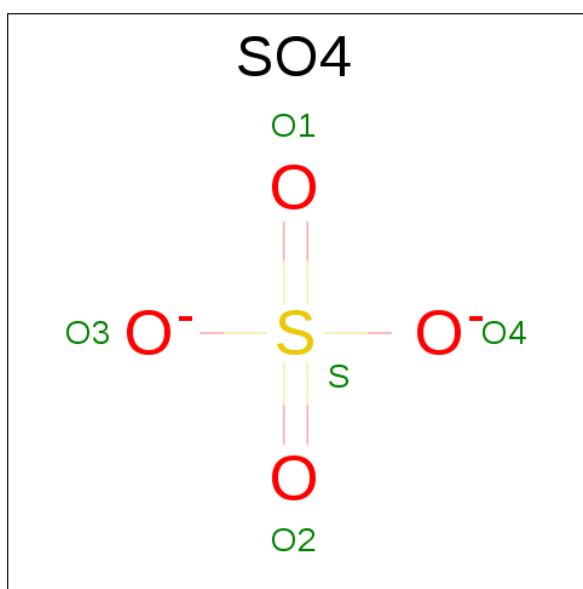
Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	MSE	-	CLONING ARTIFACT	UNP P76215
A	-3	ALA	-	CLONING ARTIFACT	UNP P76215
A	-2	GLY	-	CLONING ARTIFACT	UNP P76215
A	-1	ASP	-	CLONING ARTIFACT	UNP P76215
A	0	PRO	-	CLONING ARTIFACT	UNP P76215
A	1	MSE	MET	MODIFIED RESIDUE	UNP P76215
A	62	MSE	MET	MODIFIED RESIDUE	UNP P76215
A	102	MSE	MET	MODIFIED RESIDUE	UNP P76215
A	105	MSE	MET	MODIFIED RESIDUE	UNP P76215
A	266	MSE	MET	MODIFIED RESIDUE	UNP P76215
A	268	MSE	MET	MODIFIED RESIDUE	UNP P76215
A	276	MSE	MET	MODIFIED RESIDUE	UNP P76215
A	317	MSE	MET	MODIFIED RESIDUE	UNP P76215
A	323	LEU	-	EXPRESSION TAG	UNP P76215
A	324	GLU	-	EXPRESSION TAG	UNP P76215
A	325	HIS	-	EXPRESSION TAG	UNP P76215
A	326	HIS	-	EXPRESSION TAG	UNP P76215
A	327	HIS	-	EXPRESSION TAG	UNP P76215
A	328	HIS	-	EXPRESSION TAG	UNP P76215
A	329	HIS	-	EXPRESSION TAG	UNP P76215
A	330	HIS	-	EXPRESSION TAG	UNP P76215
B	-4	MSE	-	CLONING ARTIFACT	UNP P76215
B	-3	ALA	-	CLONING ARTIFACT	UNP P76215
B	-2	GLY	-	CLONING ARTIFACT	UNP P76215
B	-1	ASP	-	CLONING ARTIFACT	UNP P76215

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	PRO	-	CLONING ARTIFACT	UNP P76215
B	1	MSE	MET	MODIFIED RESIDUE	UNP P76215
B	62	MSE	MET	MODIFIED RESIDUE	UNP P76215
B	102	MSE	MET	MODIFIED RESIDUE	UNP P76215
B	105	MSE	MET	MODIFIED RESIDUE	UNP P76215
B	266	MSE	MET	MODIFIED RESIDUE	UNP P76215
B	268	MSE	MET	MODIFIED RESIDUE	UNP P76215
B	276	MSE	MET	MODIFIED RESIDUE	UNP P76215
B	317	MSE	MET	MODIFIED RESIDUE	UNP P76215
B	323	LEU	-	EXPRESSION TAG	UNP P76215
B	324	GLU	-	EXPRESSION TAG	UNP P76215
B	325	HIS	-	EXPRESSION TAG	UNP P76215
B	326	HIS	-	EXPRESSION TAG	UNP P76215
B	327	HIS	-	EXPRESSION TAG	UNP P76215
B	328	HIS	-	EXPRESSION TAG	UNP P76215
B	329	HIS	-	EXPRESSION TAG	UNP P76215
B	330	HIS	-	EXPRESSION TAG	UNP P76215

- 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	A	1	Total	O	S	0	0
			5	4	1		

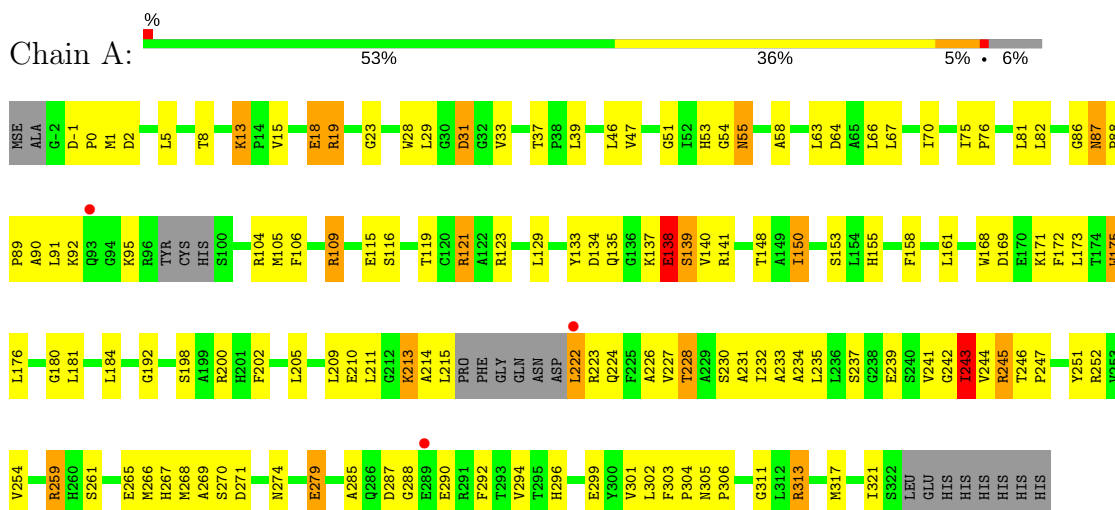
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	70	Total 70	O 70	0	0
3	B	44	Total 44	O 44	0	0

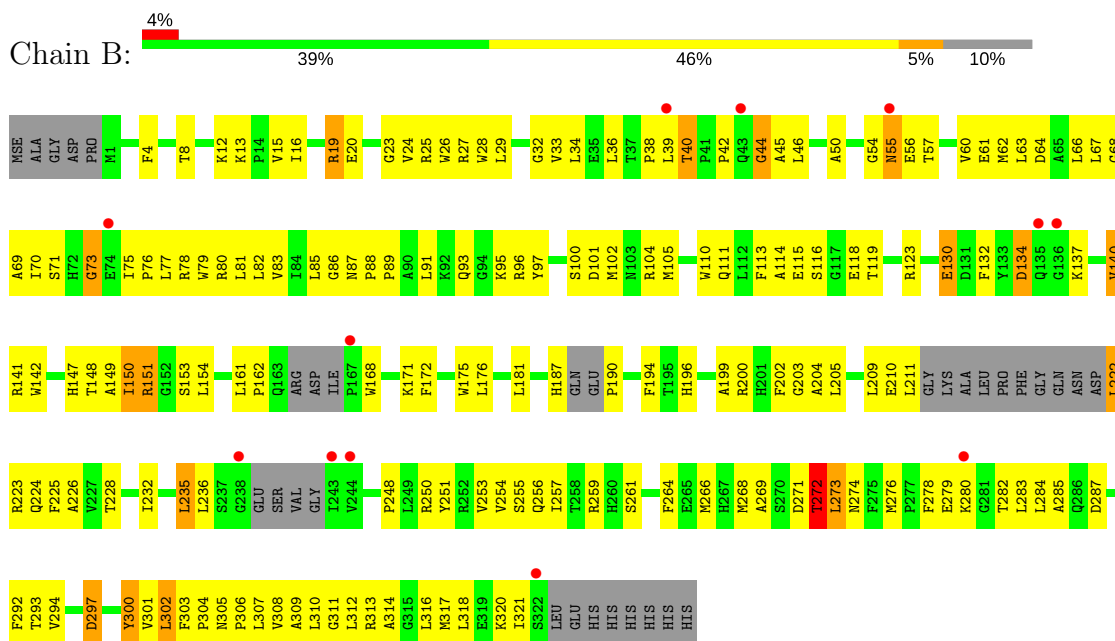
3 Residue-property plots

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: Succinylglutamate desuccinylase



• Molecule 1: Succinylglutamate desuccinylase



4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	146.80Å 146.80Å 209.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.98 – 3.10 19.98 – 3.10	Depositor EDS
% Data completeness (in resolution range)	93.4 (19.98-3.10) 97.0 (19.98-3.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.37 (at 3.09Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.227 , 0.289 0.233 , 0.284	Depositor DCC
R_{free} test set	2002 reflections (9.81%)	DCC
Wilson B-factor (Å ²)	62.1	Xtriage
Anisotropy	0.022	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 88.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.026 for -1/2*h-1/2*k-1/2*l,-1/2*h-1/2*k+1/2*l,-h+k 0.018 for -1/2*h+1/2*k-1/2*l,1/2*h-1/2*k-1/2*l,-h-k	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4969	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% 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	0.47	0/2520	0.65	0/3405
1	B	0.40	0/2429	0.56	0/3281
All	All	0.44	0/4949	0.60	0/6686

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

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	2468	0	2464	142	0
1	B	2377	0	2367	147	1
2	A	10	0	0	0	0
3	A	70	0	0	7	0
3	B	44	0	0	2	0
All	All	4969	0	4831	287	1

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 30.

All (287) 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:171:LYS:HD2	1:B:171:LYS:H	1.14	1.10
1:A:70:ILE:HG12	1:A:75:ILE:HD11	1.48	0.95
1:B:62:MSE:HE2	1:B:225:PHE:HB3	1.47	0.94
1:B:268:MSE:HE2	1:B:284:LEU:HD12	1.55	0.88
1:A:222:LEU:HG	1:A:223:ARG:H	1.42	0.85
1:A:129:LEU:HD21	1:A:198:SER:HB3	1.60	0.83
1:B:187:HIS:HB3	1:B:190:PRO:HD2	1.62	0.81
1:B:153:SER:HA	1:B:273:LEU:HB2	1.62	0.81
1:A:104:ARG:HD2	1:A:313:ARG:HH12	1.46	0.81
1:B:171:LYS:N	1:B:171:LYS:HD2	1.96	0.81
1:A:153:SER:HB3	1:A:274:ASN:HB2	1.63	0.79
1:B:116:SER:HB2	1:B:119:THR:HG23	1.65	0.79
1:A:104:ARG:HD2	1:A:313:ARG:NH1	1.98	0.77
1:A:274:ASN:HD21	1:A:303:PHE:H	1.28	0.76
1:A:245:ARG:HG3	1:A:246:THR:H	1.53	0.74
1:B:268:MSE:CE	1:B:284:LEU:HD12	2.17	0.73
1:A:55:ASN:H	1:A:55:ASN:HD22	1.37	0.73
1:B:259:ARG:HD3	1:B:309:ALA:O	1.89	0.73
1:A:104:ARG:HD3	3:A:414:HOH:O	1.90	0.71
1:B:269:ALA:O	1:B:272:THR:HG22	1.89	0.71
1:A:55:ASN:N	1:A:55:ASN:HD22	1.89	0.71
1:A:213:LYS:HB2	1:A:213:LYS:NZ	2.07	0.70
1:B:57:THR:O	1:B:60:VAL:HG12	1.93	0.68
1:A:243:ILE:HD12	1:A:247:PRO:HD3	1.75	0.68
1:A:306:PRO:HG2	3:A:410:HOH:O	1.92	0.67
1:B:301:VAL:HG22	1:B:317:MSE:SE	2.44	0.67
1:B:278:PHE:HD2	1:B:282:THR:HG21	1.58	0.67
1:B:113:PHE:O	1:B:119:THR:HG21	1.96	0.66
1:B:264:PHE:HA	1:B:287:ASP:HA	1.77	0.66
1:A:222:LEU:N	1:A:222:LEU:HD23	2.11	0.66
1:B:142:TRP:NE1	1:B:205:LEU:HD12	2.11	0.65
1:B:256:GLN:HG2	1:B:316:LEU:HD23	1.80	0.64
1:B:151:ARG:HE	1:B:304:PRO:HG2	1.63	0.63
1:B:172:PHE:HB2	1:B:205:LEU:HD13	1.81	0.63
1:A:109:ARG:HH21	1:A:109:ARG:HG2	1.64	0.63
1:A:175:TRP:HZ3	1:A:231:ALA:O	1.82	0.62
1:A:155:HIS:HE1	1:A:251:TYR:OH	1.84	0.61
1:B:101:ASP:HB3	1:B:104:ARG:HD3	1.83	0.61
1:B:259:ARG:CZ	1:B:310:LEU:HG	2.30	0.61
1:B:63:LEU:HD21	1:B:83:VAL:HG13	1.81	0.61
1:A:46:LEU:HD12	1:A:47:VAL:N	2.16	0.60
1:B:13:LYS:HG3	1:B:28:TRP:CH2	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:310:LEU:HD12	1:B:310:LEU:H	1.65	0.60
1:B:142:TRP:HE1	1:B:205:LEU:HD12	1.65	0.60
1:A:31:ASP:O	1:A:121:ARG:NH1	2.33	0.60
1:A:153:SER:CB	1:A:274:ASN:HB2	2.30	0.60
1:A:104:ARG:CZ	1:A:313:ARG:HG2	2.30	0.60
1:A:296:HIS:HB2	1:A:299:GLU:CG	2.31	0.60
1:B:45:ALA:HA	1:B:80:ARG:O	2.02	0.60
1:A:243:ILE:HG12	1:A:243:ILE:O	2.02	0.60
1:B:302:LEU:HD12	1:B:318:LEU:HD11	1.84	0.60
1:A:296:HIS:HE1	3:A:409:HOH:O	1.85	0.59
1:A:259:ARG:HG3	1:A:259:ARG:O	2.02	0.59
1:B:70:ILE:HD13	1:B:77:LEU:HD23	1.85	0.59
1:A:226:ALA:HB3	3:A:412:HOH:O	2.03	0.58
1:B:181:LEU:O	1:B:248:PRO:HB3	2.03	0.58
1:A:279:GLU:CD	1:A:279:GLU:H	2.06	0.58
1:A:243:ILE:HD11	1:A:246:THR:HA	1.86	0.58
1:B:140:VAL:O	1:B:141:ARG:HB2	2.04	0.57
1:B:67:LEU:HD21	1:B:83:VAL:HG21	1.86	0.57
1:B:153:SER:HB3	1:B:274:ASN:HB2	1.87	0.57
1:B:68:GLY:O	1:B:71:SER:HB3	2.04	0.57
1:A:296:HIS:HB2	1:A:299:GLU:HG3	1.84	0.57
1:B:142:TRP:CZ3	1:B:235:LEU:HD11	2.40	0.57
1:B:284:LEU:HD22	1:B:284:LEU:H	1.68	0.57
1:A:28:TRP:CE3	1:A:28:TRP:O	2.58	0.57
1:A:274:ASN:ND2	1:A:303:PHE:H	2.02	0.57
1:B:273:LEU:H	1:B:273:LEU:HD23	1.68	0.56
1:B:45:ALA:HA	1:B:80:ARG:HB3	1.86	0.56
1:B:268:MSE:SE	1:B:301:VAL:HG21	2.53	0.56
1:A:104:ARG:NH2	1:A:313:ARG:HG2	2.20	0.56
1:B:256:GLN:HG2	1:B:316:LEU:CD2	2.35	0.56
1:A:47:VAL:HG22	1:A:82:LEU:HD23	1.87	0.56
1:A:18:GLU:O	1:A:19:ARG:HB2	2.06	0.56
1:A:-1:ASP:HB3	1:A:2:ASP:OD2	2.05	0.56
1:B:105:MSE:HE1	1:B:118:GLU:C	2.26	0.55
1:B:266:MSE:HA	1:B:285:ALA:HB2	1.88	0.55
1:B:320:LYS:NZ	1:B:320:LYS:HB3	2.22	0.55
1:B:306:PRO:HG2	1:B:307:LEU:HD22	1.86	0.55
1:A:198:SER:HA	1:A:202:PHE:HD1	1.71	0.55
1:B:171:LYS:CD	1:B:171:LYS:H	2.01	0.55
1:A:133:TYR:CD1	1:A:141:ARG:HD2	2.41	0.55
1:A:86:GLY:HA2	1:A:121:ARG:NH2	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:LYS:HA	3:A:438:HOH:O	2.07	0.55
1:B:24:VAL:HA	1:B:38:PRO:HA	1.87	0.55
1:A:223:ARG:HH21	1:A:223:ARG:N	2.05	0.55
1:A:58:ALA:HB3	1:A:148:THR:HG21	1.88	0.55
1:A:252:ARG:HB3	1:A:321:ILE:HD11	1.89	0.54
1:B:199:ALA:HA	1:B:204:ALA:H	1.72	0.54
1:B:309:ALA:HB3	1:B:312:LEU:HD13	1.90	0.54
1:B:85:LEU:O	1:B:88:PRO:HD3	2.08	0.54
1:B:153:SER:CB	1:B:274:ASN:HB2	2.38	0.54
1:B:259:ARG:HD3	1:B:310:LEU:HA	1.89	0.54
1:B:282:THR:HB	1:B:294:VAL:HG23	1.89	0.54
1:A:181:LEU:HD11	1:A:209:LEU:HG	1.89	0.54
1:B:255:SER:HB3	1:B:317:MSE:HB2	1.90	0.54
1:B:8:THR:HG21	1:B:32:GLY:HA2	1.89	0.54
1:A:230:SER:O	1:A:233:ALA:HB3	2.07	0.53
1:A:13:LYS:HD3	1:A:13:LYS:H	1.74	0.53
1:A:47:VAL:HG22	1:A:82:LEU:HB3	1.91	0.53
1:B:257:ILE:HG13	1:B:292:PHE:CD2	2.44	0.53
1:B:310:LEU:HD12	1:B:310:LEU:N	2.23	0.53
1:B:130:GLU:HA	1:B:202:PHE:CE2	2.43	0.53
1:A:259:ARG:NH2	1:A:311:GLY:O	2.41	0.53
1:A:70:ILE:CG1	1:A:75:ILE:HD11	2.30	0.53
1:A:173:LEU:HD22	1:A:184:LEU:HD13	1.90	0.53
1:A:175:TRP:CZ3	1:A:231:ALA:O	2.62	0.52
1:B:149:ALA:H	1:B:211:LEU:C	2.12	0.52
1:B:161:LEU:HD12	1:B:161:LEU:N	2.24	0.52
1:B:40:THR:O	1:B:42:PRO:HD3	2.08	0.52
1:A:172:PHE:HB2	1:A:205:LEU:HD13	1.91	0.52
1:A:223:ARG:HH21	1:A:223:ARG:H	1.57	0.52
1:A:88:PRO:N	1:A:89:PRO:HD2	2.24	0.52
1:B:150:ILE:H	1:B:150:ILE:HD13	1.73	0.52
1:A:106:PHE:HB2	1:A:192:GLY:O	2.10	0.52
1:A:133:TYR:CE1	1:A:141:ARG:HB3	2.45	0.52
1:A:115:GLU:HG2	1:B:114:ALA:HA	1.90	0.52
1:A:175:TRP:HH2	1:A:232:ILE:HA	1.74	0.52
1:B:24:VAL:N	1:B:39:LEU:HD23	2.24	0.52
1:A:215:LEU:H	1:A:215:LEU:HD23	1.74	0.51
1:A:1:MSE:HE2	1:A:63:LEU:O	2.10	0.51
1:B:259:ARG:HG3	1:B:264:PHE:HB3	1.91	0.51
1:B:54:GLY:C	1:B:56:GLU:H	2.12	0.51
1:B:26:TRP:CE2	1:B:36:LEU:HD11	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:LYS:HB2	1:A:213:LYS:HZ3	1.75	0.51
1:B:101:ASP:HB3	1:B:104:ARG:CD	2.40	0.51
1:B:70:ILE:HD11	1:B:236:LEU:HD23	1.91	0.51
1:A:46:LEU:HD12	1:A:47:VAL:H	1.76	0.51
1:A:227:VAL:HG13	1:A:228:THR:N	2.26	0.50
1:B:100:SER:O	1:B:118:GLU:HG3	2.11	0.50
1:A:55:ASN:ND2	1:A:55:ASN:H	2.07	0.50
1:B:223:ARG:HB3	1:B:226:ALA:HB2	1.93	0.50
1:A:86:GLY:HA2	1:A:121:ARG:HH21	1.76	0.50
1:B:269:ALA:C	1:B:271:ASP:H	2.14	0.50
1:B:300:TYR:N	1:B:300:TYR:CD1	2.79	0.50
1:A:222:LEU:HG	1:A:223:ARG:N	2.21	0.50
1:A:88:PRO:HA	1:A:91:LEU:HD12	1.94	0.50
1:A:244:VAL:HG23	1:A:245:ARG:HG2	1.94	0.49
1:A:304:PRO:HB2	3:A:410:HOH:O	2.11	0.49
1:A:305:ASN:HB3	1:A:306:PRO:HD3	1.93	0.49
1:B:259:ARG:CD	1:B:310:LEU:HA	2.42	0.49
1:B:142:TRP:CE3	1:B:235:LEU:HD11	2.48	0.49
1:A:76:PRO:HG2	1:A:237:SER:CB	2.43	0.49
1:B:222:LEU:N	1:B:222:LEU:HD12	2.28	0.49
1:B:176:LEU:HB3	1:B:181:LEU:HD12	1.93	0.49
1:B:175:TRP:CE2	1:B:235:LEU:HD12	2.48	0.49
1:A:274:ASN:ND2	1:A:302:LEU:HA	2.28	0.49
1:A:234:ALA:HB1	1:A:239:GLU:HB2	1.95	0.49
1:A:67:LEU:HD13	1:A:70:ILE:HD12	1.95	0.49
1:B:61:GLU:O	1:B:64:ASP:HB3	2.13	0.49
1:A:23:GLY:C	1:A:39:LEU:HD23	2.33	0.48
1:B:302:LEU:HD22	1:B:303:PHE:CD2	2.48	0.48
1:B:25:ARG:HD3	1:B:27:ARG:HG3	1.96	0.48
1:A:222:LEU:HG	1:A:223:ARG:NH2	2.29	0.48
1:B:82:LEU:HD13	1:B:132:PHE:CD1	2.49	0.48
1:B:88:PRO:HG2	1:B:89:PRO:HD3	1.95	0.48
1:A:158:PHE:CZ	1:A:210:GLU:HG3	2.48	0.48
1:A:243:ILE:HD11	1:A:245:ARG:O	2.13	0.48
1:A:134:ASP:O	1:A:137:LYS:HG3	2.14	0.48
1:B:101:ASP:HB3	1:B:104:ARG:HG2	1.96	0.48
1:B:282:THR:HB	1:B:294:VAL:CG2	2.44	0.48
1:A:175:TRP:CH2	1:A:232:ILE:HA	2.49	0.47
1:A:88:PRO:HG2	1:A:89:PRO:HD3	1.96	0.47
1:B:257:ILE:HD12	1:B:257:ILE:N	2.29	0.47
1:B:266:MSE:HA	1:B:285:ALA:CB	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:LEU:HB3	1:A:168:TRP:CZ2	2.49	0.47
1:B:280:LYS:C	1:B:280:LYS:HD3	2.35	0.47
1:A:175:TRP:HH2	1:A:232:ILE:CA	2.27	0.47
1:B:162:PRO:CG	1:B:196:HIS:HA	2.44	0.47
1:B:23:GLY:C	1:B:39:LEU:HD23	2.34	0.47
1:B:38:PRO:HD3	1:B:77:LEU:HD12	1.94	0.47
1:B:56:GLU:HG2	1:B:148:THR:HB	1.96	0.47
1:B:302:LEU:HB3	1:B:316:LEU:HB2	1.97	0.47
1:B:141:ARG:O	1:B:142:TRP:HD1	1.98	0.47
1:B:4:PHE:O	1:B:8:THR:HG23	2.14	0.47
1:A:222:LEU:HG	1:A:223:ARG:HH22	1.79	0.47
1:A:313:ARG:O	1:A:313:ARG:HD2	2.15	0.47
1:A:175:TRP:HH2	1:A:232:ILE:N	2.12	0.47
1:B:27:ARG:O	1:B:34:LEU:HD12	2.14	0.47
1:A:105:MSE:O	1:A:123:ARG:NH2	2.48	0.46
1:B:153:SER:HB2	1:B:274:ASN:O	2.16	0.46
1:A:55:ASN:ND2	1:A:55:ASN:N	2.60	0.46
1:B:259:ARG:O	1:B:311:GLY:HA2	2.15	0.46
1:B:273:LEU:N	1:B:273:LEU:HD23	2.31	0.46
1:A:266:MSE:HE3	1:A:303:PHE:HE2	1.81	0.46
1:B:280:LYS:HD3	1:B:280:LYS:O	2.16	0.46
1:B:73:GLY:C	1:B:75:ILE:H	2.18	0.46
1:A:213:LYS:HB2	1:A:213:LYS:HZ2	1.81	0.46
1:A:243:ILE:HG23	3:A:446:HOH:O	2.15	0.46
1:B:261:SER:H	1:B:287:ASP:CG	2.20	0.46
1:A:175:TRP:CZ3	1:A:231:ALA:C	2.90	0.45
1:A:66:LEU:O	1:A:70:ILE:HG13	2.15	0.45
1:A:287:ASP:H	1:A:290:GLU:HB2	1.82	0.45
1:A:29:LEU:HB2	1:A:33:VAL:HG12	1.98	0.45
1:B:101:ASP:O	1:B:104:ARG:HG2	2.17	0.45
1:A:245:ARG:HG3	1:A:245:ARG:HH21	1.82	0.45
1:B:259:ARG:NH2	1:B:310:LEU:HG	2.31	0.45
1:A:13:LYS:H	1:A:13:LYS:CD	2.29	0.44
1:A:294:VAL:HG13	1:A:299:GLU:HG3	1.99	0.44
1:B:25:ARG:O	1:B:25:ARG:HD2	2.17	0.44
1:A:223:ARG:NH2	1:A:223:ARG:H	2.14	0.44
1:A:54:GLY:HA3	1:A:90:ALA:HB1	1.98	0.44
1:B:279:GLU:HA	1:B:297:ASP:O	2.17	0.44
1:A:87:ASN:H	1:A:121:ARG:HD2	1.81	0.44
1:B:224:GLN:H	1:B:224:GLN:HG3	1.64	0.44
1:B:29:LEU:HB2	1:B:33:VAL:HG12	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:TRP:CZ3	1:B:111:GLN:HG2	2.52	0.44
1:B:147:HIS:HB3	1:B:210:GLU:OE2	2.17	0.44
1:B:181:LEU:HD11	1:B:209:LEU:HD22	2.00	0.44
1:B:20:GLU:H	1:B:20:GLU:CD	2.21	0.44
1:A:261:SER:HB3	1:A:287:ASP:OD1	2.18	0.44
1:B:97:TYR:HB3	3:B:333:HOH:O	2.16	0.44
1:B:76:PRO:HB2	1:B:78:ARG:HH12	1.83	0.44
1:B:36:LEU:HB2	1:B:81:LEU:HB3	1.98	0.44
1:B:93:GLN:HB3	1:B:95:LYS:HG3	2.00	0.44
1:A:243:ILE:CD1	1:A:246:THR:HA	2.46	0.43
1:B:110:TRP:CH2	1:B:123:ARG:HD2	2.52	0.43
1:B:306:PRO:HG2	1:B:307:LEU:CD2	2.48	0.43
1:B:69:ALA:HB1	1:B:75:ILE:HG22	2.00	0.43
1:A:269:ALA:O	1:A:271:ASP:N	2.52	0.43
1:B:254:VAL:HG12	1:B:254:VAL:O	2.18	0.43
1:A:279:GLU:CD	1:A:279:GLU:N	2.72	0.43
1:B:268:MSE:HB3	1:B:272:THR:HG21	2.01	0.43
1:B:19:ARG:HG2	1:B:26:TRP:HB2	2.01	0.43
1:A:63:LEU:O	1:A:67:LEU:HD23	2.18	0.43
1:A:241:VAL:HG22	1:A:242:GLY:N	2.33	0.43
1:A:265:GLU:O	1:A:285:ALA:HB1	2.19	0.43
1:A:76:PRO:HG2	1:A:237:SER:HB2	2.01	0.43
1:B:116:SER:HB2	1:B:119:THR:CG2	2.44	0.43
1:B:228:THR:O	1:B:232:ILE:HG13	2.18	0.43
1:B:308:VAL:HB	1:B:314:ALA:HB2	2.01	0.43
1:A:243:ILE:CD1	1:A:247:PRO:HD3	2.48	0.42
1:A:301:VAL:HG12	1:A:302:LEU:N	2.34	0.42
1:B:175:TRP:HZ2	1:B:232:ILE:HA	1.84	0.42
1:A:1:MSE:HG2	1:A:64:ASP:HA	2.02	0.42
1:B:102:MSE:HA	1:B:105:MSE:HE2	2.01	0.42
1:A:150:ILE:HG12	1:A:150:ILE:H	1.63	0.42
1:B:305:ASN:O	1:B:308:VAL:HB	2.18	0.42
1:A:-1:ASP:OD2	1:A:0:PRO:HD2	2.20	0.42
1:B:222:LEU:C	1:B:224:GLN:H	2.23	0.42
1:B:276:MSE:O	1:B:300:TYR:HA	2.20	0.42
1:A:180:GLY:HA2	1:A:245:ARG:HH11	1.84	0.42
1:B:269:ALA:C	1:B:271:ASP:N	2.73	0.42
1:A:5:LEU:O	1:A:8:THR:HB	2.20	0.42
1:B:142:TRP:HZ3	1:B:235:LEU:HD11	1.85	0.42
1:A:313:ARG:H	1:A:313:ARG:CD	2.33	0.42
1:A:51:GLY:HA3	1:A:91:LEU:HD21	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57:THR:O	1:B:61:GLU:HG3	2.18	0.42
1:A:15:VAL:O	1:B:16:ILE:HA	2.19	0.41
1:A:211:LEU:H	1:A:211:LEU:HD22	1.85	0.41
1:A:87:ASN:H	1:A:121:ARG:HH21	1.67	0.41
1:A:215:LEU:H	1:A:215:LEU:CD2	2.34	0.41
1:B:251:TYR:N	1:B:251:TYR:CD1	2.88	0.41
1:A:294:VAL:HG13	1:A:299:GLU:CG	2.50	0.41
1:A:104:ARG:NH1	1:A:313:ARG:HH11	2.17	0.41
1:A:53:HIS:O	1:A:95:LYS:HD2	2.20	0.41
1:B:283:LEU:HA	1:B:293:THR:HG22	2.03	0.41
1:B:50:ALA:O	1:B:85:LEU:HA	2.20	0.41
1:A:169:ASP:OD2	1:A:171:LYS:HB2	2.20	0.41
1:B:15:VAL:HG23	3:B:344:HOH:O	2.20	0.41
1:B:266:MSE:HB3	1:B:306:PRO:HB3	2.02	0.41
1:B:44:GLY:HA3	1:B:79:TRP:CD2	2.55	0.41
1:A:254:VAL:HG22	1:A:317:MSE:O	2.19	0.41
1:A:274:ASN:HD21	1:A:303:PHE:N	2.07	0.41
1:A:92:LYS:O	1:A:92:LYS:HG3	2.21	0.41
1:B:222:LEU:O	1:B:224:GLN:HG3	2.20	0.41
1:B:253:VAL:HG11	1:B:316:LEU:HD22	2.03	0.41
1:A:266:MSE:C	1:A:268:MSE:H	2.23	0.41
1:A:129:LEU:HD23	1:A:202:PHE:CE1	2.55	0.41
1:B:82:LEU:HB2	1:B:132:PHE:CZ	2.56	0.41
1:A:172:PHE:CE2	1:A:176:LEU:HD11	2.56	0.41
1:A:148:THR:HG21	1:A:214:ALA:HB2	2.03	0.40
1:A:18:GLU:O	1:A:19:ARG:CB	2.69	0.40
1:A:269:ALA:C	1:A:271:ASP:H	2.25	0.40
1:B:161:LEU:HA	1:B:162:PRO:HD3	1.98	0.40
1:B:253:VAL:HA	1:B:318:LEU:HD23	2.03	0.40
1:B:54:GLY:O	1:B:55:ASN:HB2	2.21	0.40
1:B:76:PRO:HB2	1:B:78:ARG:NH1	2.36	0.40
1:A:161:LEU:HD12	1:A:161:LEU:N	2.36	0.40
1:B:272:THR:O	1:B:304:PRO:HG3	2.21	0.40
1:A:305:ASN:HB3	1:A:306:PRO:CD	2.51	0.40
1:A:116:SER:OG	1:A:119:THR:HG23	2.22	0.40
1:A:138:GLU:HB2	1:A:139:SER:H	1.54	0.40
1:A:67:LEU:HD22	1:A:67:LEU:N	2.37	0.40
1:B:118:GLU:HG2	1:B:118:GLU:H	1.70	0.40
1:A:175:TRP:CH2	1:A:231:ALA:C	2.95	0.40
1:A:81:LEU:HD12	1:A:82:LEU:H	1.86	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:LEU:CD2	1:B:154:LEU:CD2[7_555]	1.98	0.22

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	310/335 (92%)	258 (83%)	40 (13%)	12 (4%)	3	21
1	B	293/335 (88%)	236 (80%)	47 (16%)	10 (3%)	4	24
All	All	603/670 (90%)	494 (82%)	87 (14%)	22 (4%)	4	23

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	139	SER
1	A	243	ILE
1	A	19	ARG
1	A	138	GLU
1	A	245	ARG
1	A	292	PHE
1	B	86	GLY
1	B	200	ARG
1	B	203	GLY
1	A	31	ASP
1	A	109	ARG
1	A	270	SER
1	A	267	HIS
1	B	137	LYS
1	B	73	GLY
1	B	134	ASP
1	B	272	THR
1	B	321	ILE

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Mol	Chain	Res	Type
1	A	135	GLN
1	A	288	GLY
1	B	140	VAL
1	B	44	GLY

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	260/268 (97%)	240 (92%)	20 (8%)	15	48
1	B	250/268 (93%)	225 (90%)	25 (10%)	9	33
All	All	510/536 (95%)	465 (91%)	45 (9%)	12	42

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LYS
1	A	18	GLU
1	A	37	THR
1	A	55	ASN
1	A	87	ASN
1	A	121	ARG
1	A	138	GLU
1	A	140	VAL
1	A	150	ILE
1	A	175	TRP
1	A	200	ARG
1	A	213	LYS
1	A	222	LEU
1	A	224	GLN
1	A	228	THR
1	A	235	LEU
1	A	243	ILE
1	A	259	ARG
1	A	279	GLU

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Mol	Chain	Res	Type
1	A	313	ARG
1	B	12	LYS
1	B	19	ARG
1	B	40	THR
1	B	46	LEU
1	B	55	ASN
1	B	66	LEU
1	B	87	ASN
1	B	91	LEU
1	B	96	ARG
1	B	115	GLU
1	B	130	GLU
1	B	134	ASP
1	B	150	ILE
1	B	151	ARG
1	B	168	TRP
1	B	194	PHE
1	B	222	LEU
1	B	235	LEU
1	B	250	ARG
1	B	272	THR
1	B	273	LEU
1	B	297	ASP
1	B	300	TYR
1	B	302	LEU
1	B	313	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	HIS
1	A	55	ASN
1	A	72	HIS
1	A	87	ASN
1	A	155	HIS
1	A	188	GLN
1	A	267	HIS
1	A	274	ASN
1	A	286	GLN
1	A	296	HIS
1	B	22	ASN
1	B	87	ASN

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Mol	Chain	Res	Type
1	B	93	GLN
1	B	135	GLN
1	B	187	HIS
1	B	196	HIS
1	B	224	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 carbohydrates 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	$\# Z > 2$
2	SO4	A	401	-	4,4,4	0.31	0	6,6,6	0.08	0
2	SO4	A	402	-	4,4,4	0.31	0	6,6,6	0.08	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
2	SO4	A	402	-	-	0/0/0/0	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.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	308/335 (91%)	-0.38	3 (0%) 82 67	17, 42, 83, 97	0
1	B	295/335 (88%)	0.04	12 (4%) 38 18	24, 71, 97, 110	0
All	All	603/670 (90%)	-0.18	15 (2%) 58 35	17, 58, 95, 110	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	167	PRO	6.0
1	B	136	GLY	3.7
1	B	243	ILE	3.7
1	B	39	LEU	3.4
1	B	74	GLU	3.2
1	B	322	SER	3.1
1	B	135	GLN	3.0
1	B	55	ASN	3.0
1	B	43	GLN	2.5
1	A	289	GLU	2.5
1	B	280	LYS	2.4
1	A	93	GLN	2.3
1	A	222	LEU	2.3
1	B	244	VAL	2.2
1	B	238	GLY	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. 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
2	SO4	A	402	5/5	0.94	0.27	1.72	103,104,104,105	0
2	SO4	A	401	5/5	0.94	0.12	-0.77	98,98,99,99	0

6.5 Other polymers [i](#)

There are no such residues in this entry.