



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:55 AM GMT

PDB ID : 3NIV
Title : The crystal structure of Glutathione S-transferase from Legionella pneumophila
Authors : Zhang, Z.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2010-06-16
Resolution : 2.30 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

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6308 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 Glutathione S-transferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	191	Total	C	N	O	S	Se	0	1	0
			1552	1006	256	279	5	6			
1	B	188	Total	C	N	O	S	Se	0	0	0
			1525	989	250	275	5	6			
1	C	191	Total	C	N	O	S	Se	0	0	0
			1550	1004	256	279	5	6			
1	D	192	Total	C	N	O	S	Se	0	1	0
			1552	1006	256	279	5	6			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	EXPRESSION TAG	UNP Q5ZT82
A	0	SER	-	EXPRESSION TAG	UNP Q5ZT82
A	1	LEU	-	EXPRESSION TAG	UNP Q5ZT82
A	213	GLU	-	EXPRESSION TAG	UNP Q5ZT82
A	214	GLY	-	EXPRESSION TAG	UNP Q5ZT82
A	215	HIS	-	EXPRESSION TAG	UNP Q5ZT82
A	216	HIS	-	EXPRESSION TAG	UNP Q5ZT82
A	217	HIS	-	EXPRESSION TAG	UNP Q5ZT82
A	218	HIS	-	EXPRESSION TAG	UNP Q5ZT82
A	219	HIS	-	EXPRESSION TAG	UNP Q5ZT82
A	220	HIS	-	EXPRESSION TAG	UNP Q5ZT82
B	-1	MSE	-	EXPRESSION TAG	UNP Q5ZT82
B	0	SER	-	EXPRESSION TAG	UNP Q5ZT82
B	1	LEU	-	EXPRESSION TAG	UNP Q5ZT82
B	213	GLU	-	EXPRESSION TAG	UNP Q5ZT82
B	214	GLY	-	EXPRESSION TAG	UNP Q5ZT82
B	215	HIS	-	EXPRESSION TAG	UNP Q5ZT82
B	216	HIS	-	EXPRESSION TAG	UNP Q5ZT82
B	217	HIS	-	EXPRESSION TAG	UNP Q5ZT82
B	218	HIS	-	EXPRESSION TAG	UNP Q5ZT82
B	219	HIS	-	EXPRESSION TAG	UNP Q5ZT82

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	220	HIS	-	EXPRESSION TAG	UNP Q5ZT82
C	-1	MSE	-	EXPRESSION TAG	UNP Q5ZT82
C	0	SER	-	EXPRESSION TAG	UNP Q5ZT82
C	1	LEU	-	EXPRESSION TAG	UNP Q5ZT82
C	213	GLU	-	EXPRESSION TAG	UNP Q5ZT82
C	214	GLY	-	EXPRESSION TAG	UNP Q5ZT82
C	215	HIS	-	EXPRESSION TAG	UNP Q5ZT82
C	216	HIS	-	EXPRESSION TAG	UNP Q5ZT82
C	217	HIS	-	EXPRESSION TAG	UNP Q5ZT82
C	218	HIS	-	EXPRESSION TAG	UNP Q5ZT82
C	219	HIS	-	EXPRESSION TAG	UNP Q5ZT82
C	220	HIS	-	EXPRESSION TAG	UNP Q5ZT82
D	-1	MSE	-	EXPRESSION TAG	UNP Q5ZT82
D	0	SER	-	EXPRESSION TAG	UNP Q5ZT82
D	1	LEU	-	EXPRESSION TAG	UNP Q5ZT82
D	213	GLU	-	EXPRESSION TAG	UNP Q5ZT82
D	214	GLY	-	EXPRESSION TAG	UNP Q5ZT82
D	215	HIS	-	EXPRESSION TAG	UNP Q5ZT82
D	216	HIS	-	EXPRESSION TAG	UNP Q5ZT82
D	217	HIS	-	EXPRESSION TAG	UNP Q5ZT82
D	218	HIS	-	EXPRESSION TAG	UNP Q5ZT82
D	219	HIS	-	EXPRESSION TAG	UNP Q5ZT82
D	220	HIS	-	EXPRESSION TAG	UNP Q5ZT82

- Molecule 2 is water.

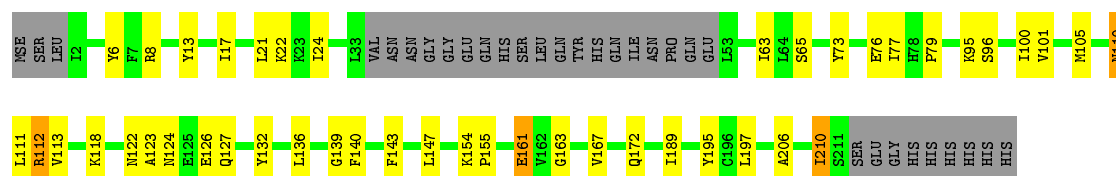
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	23	Total O 23 23	0	0
2	B	33	Total O 33 33	0	0
2	C	44	Total O 44 44	0	0
2	D	29	Total O 29 29	0	0

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 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.

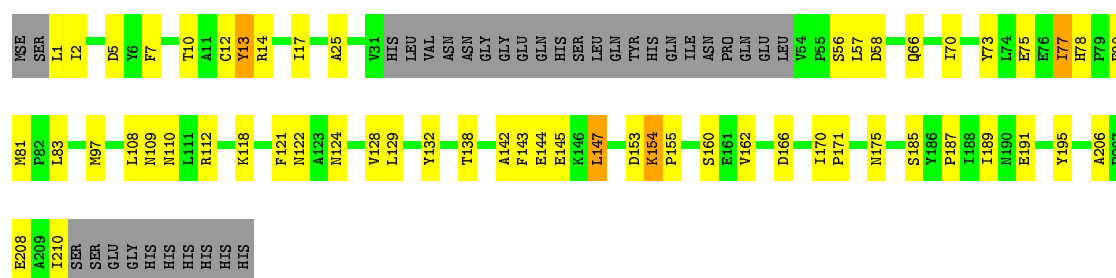
• Molecule 1: Glutathione S-transferase

Chain A: 



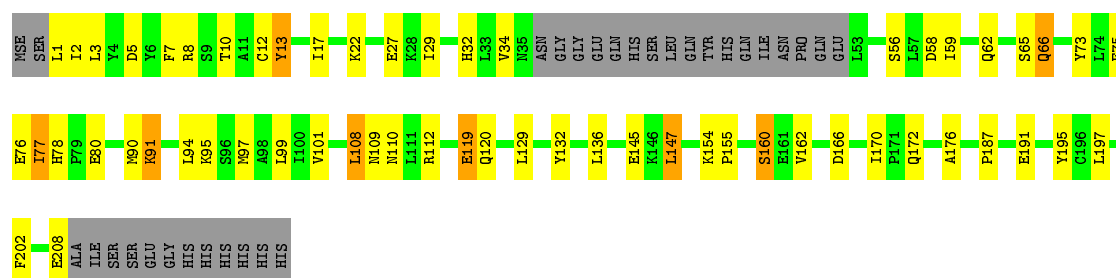
• Molecule 1: Glutathione S-transferase

Chain B: 



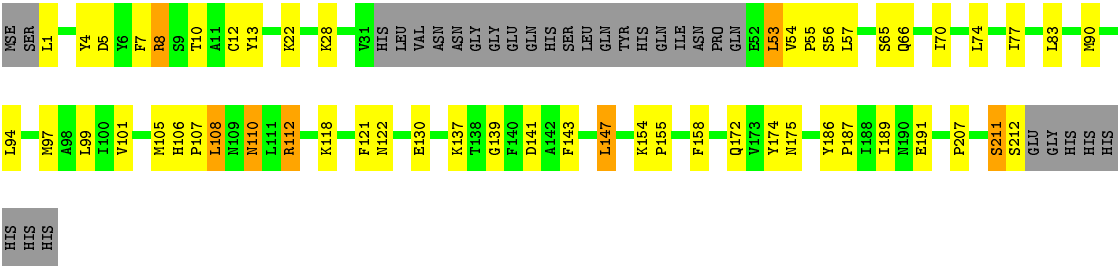
• Molecule 1: Glutathione S-transferase

Chain C: 



• Molecule 1: Glutathione S-transferase

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	51.48Å 55.04Å 84.72Å 71.99° 84.33° 68.79°	Depositor
Resolution (Å)	80.56 – 2.30 80.56 – 2.30	Depositor EDS
% Data completeness (in resolution range)	91.2 (80.56-2.30) 82.9 (80.56-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.197 , 0.249 0.201 , 0.200	Depositor DCC
R_{free} test set	3306 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	32.6	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.8	EDS
Estimated twinning fraction	0.012 for -h,-k,-l+1	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 33399 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6308	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.65% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.38	0/1589	0.53	0/2149
1	B	0.40	0/1557	0.53	0/2103
1	C	0.38	0/1583	0.51	0/2139
1	D	0.46	0/1588	0.53	0/2146
All	All	0.41	0/6317	0.53	0/8537

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	1552	0	1532	34	0
1	B	1525	0	1516	49	0
1	C	1550	0	1534	41	0
1	D	1552	0	1540	36	0
2	A	23	0	0	0	0
2	B	33	0	0	1	0
2	C	44	0	0	0	0
2	D	29	0	0	0	0
All	All	6308	0	6122	154	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (154) 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:D:118:LYS:O	1:D:122:ASN:HA	1.68	0.93
1:C:2:ILE:HG23	1:C:29:ILE:HD13	1.51	0.92
1:A:110:ASN:HD22	1:A:112:ARG:H	1.18	0.87
1:C:97:MSE:HE1	1:C:147:LEU:HD13	1.63	0.80
1:A:132:TYR:O	1:A:136:LEU:HD23	1.83	0.78
1:D:53:LEU:O	1:D:56:SER:HB2	1.84	0.78
1:C:154:LYS:HE3	1:C:160:SER:O	1.84	0.77
1:A:110:ASN:ND2	1:A:112:ARG:H	1.83	0.77
1:C:110:ASN:ND2	1:C:112:ARG:H	1.84	0.76
1:B:73:TYR:CZ	1:B:77:ILE:HD12	2.21	0.76
1:D:4:TYR:HB2	1:D:56:SER:HB3	1.67	0.75
1:C:2:ILE:HB	1:C:58:ASP:HB3	1.71	0.73
1:D:57:LEU:HD22	1:D:70:ILE:HG23	1.73	0.69
1:C:119:GLU:HG3	1:C:120:GLN:N	2.07	0.68
1:C:110:ASN:HD22	1:C:112:ARG:H	1.45	0.65
1:B:7:PHE:HB2	1:B:210:ILE:HB	1.77	0.65
1:D:8:ARG:NH1	1:D:174:TYR:OH	2.29	0.65
1:B:5:ASP:OD1	1:B:12:CYS:HB3	1.97	0.64
1:B:110:ASN:HD22	1:B:112:ARG:H	1.45	0.64
1:C:109:ASN:HB3	1:C:132:TYR:CE2	2.34	0.62
1:D:118:LYS:O	1:D:122:ASN:CA	2.46	0.61
1:A:110:ASN:HD22	1:A:112:ARG:N	1.94	0.61
1:D:187:PRO:O	1:D:191:GLU:HG3	2.01	0.61
1:A:105:MSE:SE	1:A:139:GLY:HA3	2.52	0.59
1:C:95:LYS:O	1:C:99:LEU:HG	2.01	0.59
1:B:142:ALA:O	1:B:145:GLU:HB3	2.03	0.59
1:B:78:HIS:NE2	1:B:81:MSE:HE2	2.18	0.58
1:B:208:GLU:HB2	2:B:242:HOH:O	2.04	0.58
1:B:97:MSE:CE	1:B:147:LEU:HD13	2.34	0.57
1:B:187:PRO:O	1:B:191:GLU:HG3	2.05	0.56
1:A:136:LEU:HD12	1:A:140:PHE:CE2	2.40	0.56
1:D:110:ASN:HD22	1:D:112:ARG:H	1.54	0.56
1:B:78:HIS:NE2	1:B:81:MSE:CE	2.68	0.55
1:B:155:PRO:HB2	1:B:195:TYR:CB	2.37	0.55
1:A:105:MSE:CE	1:A:136:LEU:HD13	2.36	0.54
1:B:153:ASP:HB2	1:B:160:SER:OG	2.08	0.54
1:D:22:LYS:HE2	1:D:83:LEU:HG	1.90	0.54
1:D:77:ILE:O	1:D:77:ILE:HG12	2.07	0.53
1:A:123:ALA:HA	1:A:127:GLN:OE1	2.08	0.53
1:C:13:TYR:O	1:C:17:ILE:HG13	2.09	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:118:LYS:O	1:B:122:ASN:HA	2.08	0.53
1:C:22:LYS:HE3	1:C:75:GLU:HA	1.90	0.53
1:A:112:ARG:HG2	1:A:113:VAL:N	2.24	0.53
1:D:90:MSE:HE3	1:D:94:LEU:CD1	2.39	0.52
1:B:154:LYS:HZ2	1:B:154:LYS:HB3	1.74	0.52
1:D:97:MSE:CE	1:D:147:LEU:HD13	2.39	0.52
1:A:124:ASN:OD1	1:A:126:GLU:HG2	2.10	0.52
1:D:137:LYS:NZ	1:D:141:ASP:OD2	2.41	0.52
1:A:22:LYS:CB	1:A:24:ILE:HD11	2.40	0.52
1:D:5:ASP:OD1	1:D:12:CYS:HB3	2.10	0.51
1:C:154:LYS:HB2	1:C:154:LYS:NZ	2.24	0.51
1:A:124:ASN:OD1	1:A:127:GLN:HG3	2.10	0.51
1:A:6:TYR:CE2	1:A:8:ARG:HB2	2.45	0.51
1:D:105:MSE:SE	1:D:139:GLY:HA3	2.61	0.51
1:D:65:SER:O	1:D:66:GLN:HB2	2.10	0.51
1:C:65:SER:O	1:C:66:GLN:HB3	2.12	0.50
1:D:53:LEU:O	1:D:56:SER:CB	2.56	0.50
1:B:97:MSE:HE1	1:B:147:LEU:HD13	1.92	0.50
1:D:143:PHE:HD1	1:D:147:LEU:HD22	1.77	0.50
1:B:143:PHE:CD1	1:B:147:LEU:HD22	2.47	0.50
1:D:90:MSE:HE2	1:D:158:PHE:CD1	2.47	0.49
1:C:65:SER:O	1:C:66:GLN:CB	2.60	0.49
1:D:101:VAL:HG13	1:D:172:GLN:HE21	1.78	0.49
1:B:1:LEU:HD21	1:B:57:LEU:HD11	1.95	0.48
1:B:155:PRO:HB2	1:B:195:TYR:HB2	1.96	0.48
1:C:73:TYR:CZ	1:C:77:ILE:HD12	2.48	0.48
1:B:73:TYR:CZ	1:B:77:ILE:CD1	2.95	0.48
1:C:155:PRO:HG2	1:C:195:TYR:CG	2.48	0.48
1:A:13:TYR:O	1:A:17:ILE:HG13	2.13	0.47
1:A:118:LYS:O	1:A:122:ASN:HA	2.14	0.47
1:A:96:SER:HB2	1:B:66:GLN:HB2	1.96	0.47
1:D:121:PHE:O	1:D:122:ASN:C	2.52	0.47
1:D:154:LYS:HB2	1:D:155:PRO:CD	2.45	0.47
1:B:13:TYR:CE1	1:B:17:ILE:HD11	2.50	0.47
1:B:1:LEU:HD23	1:B:1:LEU:C	2.35	0.47
1:C:59:ILE:O	1:C:62:GLN:HG3	2.15	0.47
1:B:162:VAL:HG13	1:B:166:ASP:HB2	1.97	0.47
1:A:197:LEU:HD23	1:A:197:LEU:HA	1.69	0.47
1:C:78:HIS:NE2	1:C:80:GLU:HB2	2.29	0.47
1:C:90:MSE:HE3	1:C:94:LEU:HD13	1.96	0.46
1:D:211:SER:O	1:D:212:SER:C	2.53	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:106:HIS:N	1:D:107:PRO:CD	2.79	0.46
1:C:109:ASN:OD1	1:C:132:TYR:HA	2.16	0.46
1:D:186:TYR:HB3	1:D:189:ILE:HD12	1.98	0.46
1:B:143:PHE:HD1	1:B:147:LEU:HD22	1.81	0.46
1:B:75:GLU:CD	1:B:83:LEU:H	2.20	0.45
1:D:10:THR:OG1	1:D:175:ASN:ND2	2.48	0.45
1:A:206:ALA:O	1:A:210:ILE:HD13	2.16	0.45
1:B:144:GLU:OE1	1:B:189:ILE:HG13	2.16	0.45
1:C:3:LEU:HD12	1:C:56:SER:O	2.17	0.45
1:A:73:TYR:CE2	1:A:77:ILE:HD13	2.51	0.45
1:C:154:LYS:HZ3	1:C:154:LYS:HB2	1.81	0.45
1:B:185:SER:C	1:B:187:PRO:HD3	2.37	0.45
1:A:147:LEU:CD2	1:A:189:ILE:HG13	2.47	0.45
1:B:145:GLU:OE1	1:B:145:GLU:HA	2.15	0.45
1:B:80:GLU:O	1:B:81:MSE:HB2	2.15	0.45
1:D:97:MSE:HE3	1:D:147:LEU:HD13	1.99	0.45
1:B:25:ALA:HA	1:C:34:VAL:O	2.17	0.45
1:A:112:ARG:NH2	1:B:138:THR:HG21	2.32	0.45
1:B:109:ASN:OD1	1:B:132:TYR:HA	2.17	0.45
1:C:136:LEU:HD11	1:C:176:ALA:HA	1.99	0.45
1:D:7:PHE:CE1	1:D:8:ARG:HD3	2.51	0.44
1:B:10:THR:O	1:B:14:ARG:HG3	2.18	0.44
1:D:174:TYR:CZ	1:D:207:PRO:HG2	2.52	0.44
1:A:161:GLU:OE1	1:A:161:GLU:HA	2.17	0.44
1:A:77:ILE:C	1:A:79:PRO:HD3	2.38	0.44
1:C:17:ILE:HD13	1:C:202:PHE:CD2	2.53	0.44
1:B:154:LYS:NZ	1:B:154:LYS:HB3	2.31	0.43
1:B:121:PHE:O	1:B:122:ASN:C	2.56	0.43
1:D:54:VAL:HA	1:D:55:PRO:C	2.39	0.43
1:C:162:VAL:HG13	1:C:166:ASP:HB2	2.00	0.43
1:A:101:VAL:HG13	1:A:172:GLN:HE21	1.82	0.43
1:C:91:LYS:HE2	1:C:91:LYS:HB2	1.64	0.43
1:C:187:PRO:O	1:C:191:GLU:HG3	2.18	0.43
1:D:90:MSE:HE3	1:D:94:LEU:HD13	2.00	0.43
1:C:7:PHE:CE1	1:C:8:ARG:HB2	2.53	0.43
1:B:124:ASN:O	1:B:128:VAL:HG23	2.18	0.43
1:C:95:LYS:HE3	1:C:95:LYS:HB2	1.77	0.43
1:C:7:PHE:CE1	1:C:8:ARG:HD2	2.54	0.43
1:D:70:ILE:O	1:D:74:LEU:HG	2.18	0.43
1:B:108:LEU:N	1:B:108:LEU:HD22	2.34	0.43
1:C:5:ASP:OD1	1:C:12:CYS:HB3	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:SER:HB2	1:B:66:GLN:CG	2.49	0.43
1:C:208:GLU:N	1:C:208:GLU:OE1	2.50	0.43
1:B:73:TYR:CE2	1:B:77:ILE:HD12	2.53	0.43
1:A:21:LEU:HD23	1:A:21:LEU:HA	1.80	0.42
1:A:100:ILE:O	1:A:105:MSE:HG3	2.19	0.42
1:D:90:MSE:HE2	1:D:158:PHE:CE1	2.54	0.42
1:B:162:VAL:HG13	1:B:166:ASP:CB	2.49	0.42
1:C:129:LEU:HD12	1:C:129:LEU:HA	1.82	0.42
1:C:154:LYS:HZ3	1:C:154:LYS:CB	2.33	0.42
1:B:70:ILE:O	1:B:73:TYR:HB3	2.20	0.42
1:B:170:ILE:HB	1:B:171:PRO:HD3	2.02	0.42
1:B:10:THR:HG21	1:B:175:ASN:ND2	2.35	0.42
1:A:22:LYS:C	1:A:24:ILE:HD12	2.40	0.42
1:A:163:GLY:O	1:A:167:VAL:HG23	2.21	0.41
1:C:101:VAL:HG13	1:C:172:GLN:HE21	1.85	0.41
1:A:110:ASN:ND2	1:A:111:LEU:N	2.68	0.41
1:A:22:LYS:CB	1:A:24:ILE:CD1	2.97	0.41
1:C:110:ASN:HD22	1:C:112:ARG:N	2.13	0.41
1:D:186:TYR:N	1:D:187:PRO:HD3	2.36	0.41
1:C:108:LEU:CD1	1:D:108:LEU:HD12	2.51	0.41
1:B:206:ALA:O	1:B:210:ILE:HG12	2.20	0.41
1:B:78:HIS:NE2	1:B:81:MSE:HE1	2.36	0.41
1:C:66:GLN:HG2	1:D:99:LEU:HB2	2.03	0.41
1:B:2:ILE:HB	1:B:58:ASP:HB3	2.02	0.41
1:C:197:LEU:HD23	1:C:197:LEU:HA	1.79	0.41
1:C:119:GLU:CG	1:C:120:GLN:N	2.81	0.41
1:B:154:LYS:HB2	1:B:155:PRO:CD	2.52	0.40
1:A:124:ASN:CG	1:A:127:GLN:HG3	2.41	0.40
1:B:80:GLU:CD	1:B:80:GLU:H	2.24	0.40
1:A:143:PHE:O	1:A:147:LEU:HB2	2.21	0.40
1:B:57:LEU:HD12	1:B:57:LEU:HA	1.94	0.40
1:A:155:PRO:HB2	1:A:195:TYR:CB	2.52	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	188/222 (85%)	179 (95%)	9 (5%)	0	100	100
1	B	184/222 (83%)	180 (98%)	4 (2%)	0	100	100
1	C	187/222 (84%)	184 (98%)	2 (1%)	1 (0%)	34	41
1	D	189/222 (85%)	183 (97%)	6 (3%)	0	100	100
All	All	748/888 (84%)	726 (97%)	21 (3%)	1 (0%)	56	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	66	GLN

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	169/190 (89%)	160 (95%)	9 (5%)	28	37
1	B	166/190 (87%)	160 (96%)	6 (4%)	42	57
1	C	169/190 (89%)	155 (92%)	14 (8%)	14	17
1	D	169/190 (89%)	158 (94%)	11 (6%)	21	27
All	All	673/760 (89%)	633 (94%)	40 (6%)	24	32

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	ILE
1	A	65	SER
1	A	76	GLU
1	A	95	LYS
1	A	110	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	112	ARG
1	A	154	LYS
1	A	161	GLU
1	A	210	ILE
1	B	13	TYR
1	B	56	SER
1	B	77	ILE
1	B	129	LEU
1	B	147	LEU
1	B	154	LYS
1	C	1	LEU
1	C	10	THR
1	C	13	TYR
1	C	27	GLU
1	C	32	HIS
1	C	76	GLU
1	C	77	ILE
1	C	91	LYS
1	C	108	LEU
1	C	119	GLU
1	C	145	GLU
1	C	147	LEU
1	C	160	SER
1	C	170	ILE
1	D	1	LEU
1	D	8	ARG
1	D	13	TYR
1	D	28	LYS
1	D	53	LEU
1	D	108	LEU
1	D	110	ASN
1	D	112	ARG
1	D	130	GLU
1	D	147	LEU
1	D	211	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	ASN
1	A	32	HIS
1	A	110	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	177	HIS
1	B	20	ASN
1	B	110	ASN
1	B	172	GLN
1	B	177	HIS
1	C	20	ASN
1	C	32	HIS
1	C	35	ASN
1	C	110	ASN
1	C	172	GLN
1	D	20	ASN
1	D	60	ASN
1	D	110	ASN
1	D	134	HIS
1	D	172	GLN
1	D	175	ASN

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](#)

There are no ligands in this entry.

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 [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	185/222 (83%)	-0.15	0 100 100	24, 35, 49, 61	0
1	B	182/222 (81%)	-0.16	0 100 100	22, 33, 48, 63	0
1	C	185/222 (83%)	-0.23	0 100 100	24, 33, 47, 55	0
1	D	186/222 (83%)	-0.18	0 100 100	21, 33, 46, 57	0
All	All	738/888 (83%)	-0.18	0 100 100	21, 34, 48, 63	0

There are no RSRZ outliers to report.

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.