Full wwPDB X-ray Structure Validation Report ⓘ

Nov 16, 2024 – 04:36 PM EST

PDB ID	:	3PNT
Title	:	Crystal Structure of the Streptococcus pyogenes NAD+ glycohydrolase SPN in complex with IFS, the Immunity Factor for SPN
Authors	:	Smith, C.L.; Stine Elam, J.; Ellenberger, T.; Ghosh, J.; Pinkner, J.S.; Hultgren, S.J.; Caparon, M.G.
Deposited on	:	2010-11-19
Resolution	:	2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references ⓘ) were used in the production of this report:

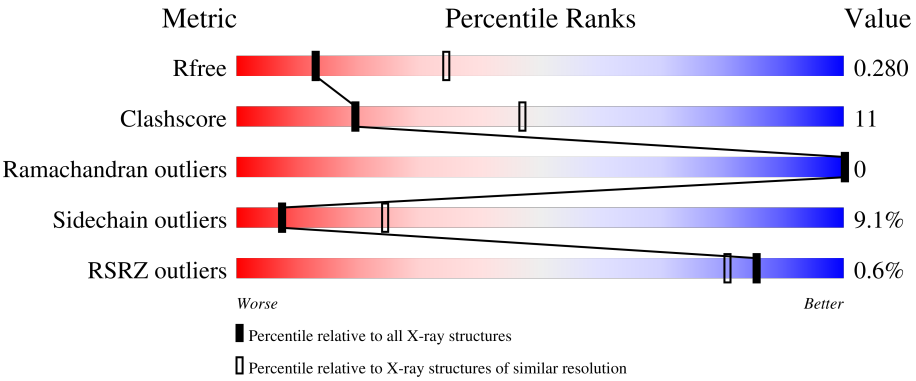
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

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 2.80 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}	164625	3657 (2.80-2.80)
Clashscore	180529	4123 (2.80-2.80)
Ramachandran outliers	177936	4071 (2.80-2.80)
Sidechain outliers	177891	4073 (2.80-2.80)
RSRZ outliers	164620	3659 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	268	73% 17% • 7%
1	C	268	% 78% 13% • 7%
2	B	161	71% 25% •
2	D	161	69% 25% 6% •

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6793 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+-glycohydrolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	250	Total	C	N	O	S	Se	0	0	0
			2001	1268	333	394	1	5			
1	C	250	Total	C	N	O	S	Se	0	0	0
			2001	1268	333	394	1	5			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	184	MSE	-	expression tag	UNP D7S065
A	185	HIS	-	expression tag	UNP D7S065
A	186	HIS	-	expression tag	UNP D7S065
A	187	HIS	-	expression tag	UNP D7S065
A	188	HIS	-	expression tag	UNP D7S065
A	189	HIS	-	expression tag	UNP D7S065
A	190	HIS	-	expression tag	UNP D7S065
C	184	MSE	-	expression tag	UNP D7S065
C	185	HIS	-	expression tag	UNP D7S065
C	186	HIS	-	expression tag	UNP D7S065
C	187	HIS	-	expression tag	UNP D7S065
C	188	HIS	-	expression tag	UNP D7S065
C	189	HIS	-	expression tag	UNP D7S065
C	190	HIS	-	expression tag	UNP D7S065

- Molecule 2 is a protein called Immunity factor for SPN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	161	Total	C	N	O	Se	0	0	0
			1323	847	215	253	8			
2	D	160	Total	C	N	O	Se	0	0	0
			1315	842	214	252	7			

- Molecule 3 is water.

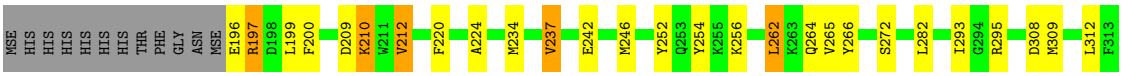
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	37	Total O 37 37	0	0
3	B	37	Total O 37 37	0	0
3	C	39	Total O 39 39	0	0
3	D	40	Total O 40 40	0	0

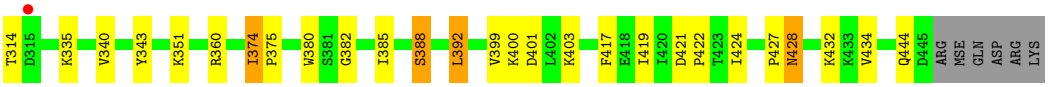
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: NAD+-glycohydrolase

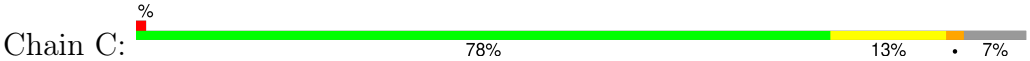
Chain A:

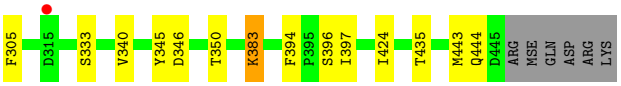




• Molecule 1: NAD+-glycohydrolase

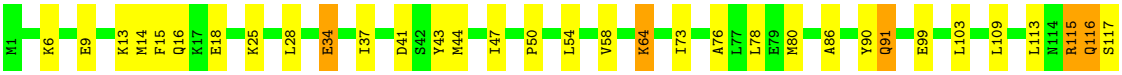
Chain C:





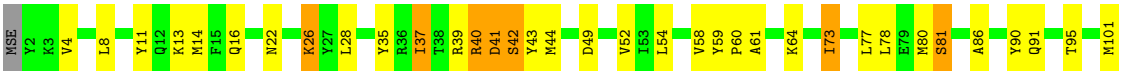
• Molecule 2: Immunity factor for SPN

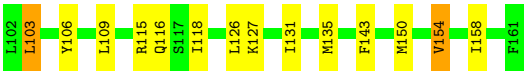
Chain B:



• Molecule 2: Immunity factor for SPN

Chain D:




L102
L103
Y106
L109
R115
Q116
S117
I118
L126
K127
I131
M135
F143
M150
V154
I158
F161

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	198.82Å 57.89Å 89.70Å 90.00° 107.20° 90.00°	Depositor
Resolution (Å)	42.72 – 2.80 42.72 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (42.72-2.80) 99.7 (42.72-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.73 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5.0093	Depositor
R, R_{free}	0.208 , 0.278 0.216 , 0.280	Depositor DCC
R_{free} test set	1240 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å^2)	52.7	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 19.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.52$, $< L^2 > = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	6793	wwPDB-VP
Average B, all atoms (Å^2)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% of the height of the origin peak. No significant pseudotranslation is detected.*

[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

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	0.57	0/2031	0.62	0/2713
1	C	0.66	1/2031 (0.0%)	0.67	2/2713 (0.1%)
2	B	0.53	0/1342	0.66	1/1796 (0.1%)
2	D	0.69	0/1334	0.65	0/1786
All	All	0.61	1/6738 (0.0%)	0.65	3/9008 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	247	TYR	CD2-CE2	-5.05	1.31	1.39

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	294	GLY	N-CA-C	9.59	137.06	113.10
1	C	295	ARG	N-CA-C	5.38	125.53	111.00
2	B	156	ARG	NE-CZ-NH1	5.14	122.87	120.30

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	2001	0	1993	33	0
1	C	2001	0	1993	29	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1323	0	1310	52	0
2	D	1315	0	1298	42	1
3	A	37	0	0	4	0
3	B	37	0	0	4	0
3	C	39	0	0	4	0
3	D	40	0	0	3	0
All	All	6793	0	6594	151	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 11.

All (151) 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:234:MSE:HE1	3:A:167:HOH:O	1.62	0.97
2:B:135:MSE:HG2	2:B:150:MSE:HE2	1.52	0.90
1:A:400:LYS:HE2	1:A:401:ASP:OD2	1.72	0.90
1:C:293:ILE:O	1:C:396:SER:HB3	1.72	0.90
2:B:90:TYR:HB2	2:B:150:MSE:HE3	1.56	0.86
2:D:44:MSE:CE	3:D:180:HOH:O	2.25	0.84
2:B:86:ALA:O	2:B:150:MSE:HE1	1.80	0.82
2:B:86:ALA:C	2:B:150:MSE:HE1	2.00	0.81
1:A:400:LYS:CE	1:A:401:ASP:OD2	2.30	0.80
1:C:246:MSE:HE1	1:C:254:TYR:CE1	2.23	0.74
2:D:135:MSE:HG2	2:D:150:MSE:HE2	1.69	0.74
1:A:209:ASP:O	1:A:212:VAL:HG13	1.87	0.73
2:B:44:MSE:HE1	2:B:143:PHE:CD1	2.24	0.73
2:D:39:ARG:HB3	2:D:52:VAL:HG21	1.70	0.73
1:C:294:GLY:O	1:C:295:ARG:CB	2.30	0.71
2:D:44:MSE:HE3	3:D:180:HOH:O	1.87	0.71
2:B:14:MSE:HE1	2:B:37:ILE:HD11	1.72	0.70
2:B:116:GLN:HG3	2:B:117:SER:H	1.54	0.70
2:B:47:ILE:HD12	2:B:91:GLN:HE22	1.57	0.70
2:B:13:LYS:HD2	2:B:13:LYS:N	2.07	0.69
2:B:14:MSE:HE3	2:B:15:PHE:CE2	2.28	0.69
1:C:259:VAL:HG12	1:C:263:LYS:HE2	1.75	0.69
1:C:211:TRP:CH2	2:D:40:ARG:HG2	2.28	0.68
2:D:86:ALA:C	2:D:150:MSE:HE1	2.14	0.68
1:C:246:MSE:HE3	1:C:252:TYR:HB3	1.76	0.68
2:B:116:GLN:HG3	2:B:117:SER:N	2.09	0.68
1:A:400:LYS:O	1:A:401:ASP:HB2	1.92	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:14:MSE:CE	2:B:15:PHE:CE2	2.77	0.67
1:A:246:MSE:HE1	1:A:254:TYR:CE2	2.29	0.67
1:A:427:PRO:O	3:A:165:HOH:O	2.13	0.66
2:D:86:ALA:HB1	2:D:150:MSE:HE1	1.78	0.65
2:D:41:ASP:N	2:D:41:ASP:OD1	2.29	0.64
1:A:428:ASN:H	1:A:428:ASN:HD22	1.43	0.64
2:D:39:ARG:NH1	2:D:41:ASP:OD2	2.30	0.64
2:B:44:MSE:HE2	3:B:188:HOH:O	1.98	0.63
2:B:76:ALA:C	2:B:80:MSE:HE3	2.19	0.63
2:B:14:MSE:HE1	2:B:37:ILE:CD1	2.29	0.63
2:D:44:MSE:HE1	2:D:143:PHE:CE1	2.34	0.62
1:A:399:VAL:HG23	1:A:399:VAL:O	1.99	0.62
1:C:234:MSE:O	1:C:235:SER:HB2	2.00	0.62
1:C:340:VAL:HG13	1:C:443:MSE:HB2	1.82	0.62
1:C:294:GLY:O	1:C:299:GLY:O	2.18	0.61
1:C:244:TYR:O	1:C:248:VAL:HG23	2.00	0.60
2:B:14:MSE:CE	2:B:15:PHE:HE2	2.15	0.60
2:D:103:LEU:O	2:D:106:TYR:O	2.20	0.60
2:B:44:MSE:CE	3:B:188:HOH:O	2.50	0.59
2:D:115:ARG:HA	2:D:118:ILE:HD12	1.85	0.59
2:B:118:ILE:HD12	2:B:118:ILE:N	2.17	0.59
1:C:294:GLY:O	1:C:295:ARG:HB3	2.02	0.59
2:B:28:LEU:O	2:B:91:GLN:HG2	2.04	0.58
1:C:340:VAL:HG13	1:C:340:VAL:O	2.04	0.58
2:D:77:LEU:HD11	2:D:95:THR:HG21	1.84	0.58
1:C:196:GLU:N	3:C:131:HOH:O	2.36	0.58
2:D:86:ALA:HB1	2:D:150:MSE:CE	2.35	0.57
1:C:305:PHE:N	3:C:42:HOH:O	2.37	0.56
2:B:135:MSE:CG	2:B:150:MSE:HE2	2.31	0.55
2:D:135:MSE:CG	2:D:150:MSE:HE2	2.37	0.55
1:A:246:MSE:HE3	1:A:252:TYR:CA	2.38	0.54
1:A:246:MSE:HE3	1:A:252:TYR:HA	1.90	0.54
2:D:44:MSE:HE1	3:D:180:HOH:O	1.97	0.54
1:A:272:SER:OG	1:A:308:ASP:OD2	2.10	0.53
1:C:269:SER:HB2	3:C:42:HOH:O	2.07	0.53
2:D:131:ILE:HG22	2:D:135:MSE:HE2	1.89	0.53
1:A:246:MSE:HE2	1:A:375:PRO:HB3	1.91	0.53
2:B:154:VAL:O	2:B:158:ILE:HD13	2.09	0.53
1:C:294:GLY:O	1:C:295:ARG:HB2	2.08	0.53
1:C:397:ILE:HD13	1:C:424:ILE:HG22	1.90	0.53
2:B:150:MSE:O	2:B:154:VAL:HG23	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:26:LYS:NZ	2:D:35:TYR:CZ	2.77	0.53
1:C:295:ARG:NH2	2:D:43:TYR:O	2.42	0.52
2:D:78:LEU:HD23	2:D:78:LEU:O	2.09	0.52
2:B:44:MSE:HE1	2:B:143:PHE:CE1	2.44	0.51
1:A:295:ARG:NH2	2:B:43:TYR:O	2.43	0.51
2:B:58:VAL:HG21	2:B:73:ILE:HD11	1.92	0.51
2:D:126:LEU:HG	2:D:135:MSE:HE1	1.91	0.51
2:D:154:VAL:HG23	2:D:158:ILE:HD12	1.93	0.51
2:B:118:ILE:N	2:B:118:ILE:CD1	2.74	0.50
1:A:400:LYS:HE3	1:A:401:ASP:OD2	2.07	0.50
2:B:25:LYS:HA	2:B:80:MSE:HE1	1.93	0.50
2:B:16:GLN:OE1	2:B:64:LYS:NZ	2.44	0.50
1:A:428:ASN:HD22	1:A:428:ASN:N	2.01	0.50
2:B:47:ILE:HD12	2:B:91:GLN:NE2	2.24	0.50
1:C:197:ARG:HA	1:C:200:PHE:HB3	1.94	0.50
1:A:308:ASP:O	1:A:312:LEU:HB2	2.12	0.49
2:D:42:SER:OG	2:D:49:ASP:OD2	2.22	0.49
1:A:385:ILE:O	1:A:388:SER:HB2	2.13	0.49
2:B:76:ALA:O	2:B:80:MSE:HE3	2.13	0.48
1:C:246:MSE:HE3	1:C:252:TYR:CB	2.42	0.48
2:B:86:ALA:O	2:B:150:MSE:CE	2.55	0.48
2:B:115:ARG:HD3	3:B:184:HOH:O	2.12	0.48
1:A:220:PHE:CZ	1:A:374:ILE:HD11	2.48	0.48
1:A:293:ILE:HD13	1:A:343:TYR:CE1	2.48	0.47
2:B:14:MSE:HE2	2:B:15:PHE:HE2	1.79	0.47
1:C:211:TRP:CZ2	2:D:40:ARG:HG2	2.48	0.47
2:D:90:TYR:HB2	2:D:150:MSE:HE3	1.96	0.47
2:B:50:PRO:O	2:B:54:LEU:HD13	2.15	0.47
2:B:90:TYR:HB2	2:B:150:MSE:CE	2.36	0.47
2:D:101:MSE:HA	2:D:101:MSE:HE2	1.96	0.47
2:D:77:LEU:O	2:D:81:SER:HB2	2.14	0.47
2:B:118:ILE:CD1	2:B:118:ILE:H	2.28	0.47
2:D:59:TYR:HB2	2:D:60:PRO:HD3	1.96	0.47
1:C:234:MSE:O	1:C:235:SER:CB	2.63	0.46
2:B:86:ALA:HB1	2:B:150:MSE:CE	2.45	0.46
1:C:230:LYS:O	1:C:234:MSE:HG3	2.16	0.46
2:B:41:ASP:OD1	2:B:41:ASP:N	2.43	0.46
2:B:139:LYS:NZ	2:B:146:TYR:O	2.48	0.46
1:A:424:ILE:HD13	1:A:434:VAL:HG21	1.98	0.46
2:D:86:ALA:CB	2:D:150:MSE:HE1	2.46	0.46
1:A:262:LEU:O	1:A:265:VAL:HG12	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:86:ALA:CA	2:B:150:MSE:HE1	2.45	0.46
2:B:135:MSE:HE2	2:B:150:MSE:CE	2.45	0.46
2:D:126:LEU:CD1	2:D:135:MSE:HE1	2.46	0.45
2:D:14:MSE:SE	2:D:37:ILE:HD11	2.67	0.45
2:D:28:LEU:O	2:D:91:GLN:HG2	2.16	0.45
1:A:246:MSE:HE3	1:A:252:TYR:HB3	1.98	0.45
1:C:274:VAL:HG13	1:C:340:VAL:HG23	1.98	0.45
1:A:419:ILE:HD13	3:A:89:HOH:O	2.16	0.45
2:B:86:ALA:HB1	2:B:150:MSE:HE1	1.98	0.45
2:B:118:ILE:HG22	2:B:122:ILE:CD1	2.47	0.44
1:A:421:ASP:N	1:A:422:PRO:HD3	2.32	0.44
2:B:119:LEU:HA	2:B:122:ILE:HD12	2.00	0.44
2:D:126:LEU:HD21	2:D:154:VAL:CG2	2.47	0.44
2:B:119:LEU:HD11	2:B:161:PHE:CD2	2.53	0.44
1:A:385:ILE:CG2	3:A:68:HOH:O	2.67	0.43
1:C:246:MSE:HE3	1:C:252:TYR:CA	2.48	0.43
2:D:8:LEU:HD13	2:D:11:TYR:HD1	1.83	0.43
2:D:58:VAL:HG21	2:D:73:ILE:CD1	2.49	0.43
2:D:39:ARG:NH1	2:D:41:ASP:CG	2.72	0.43
2:D:16:GLN:HG2	2:D:61:ALA:HA	2.00	0.43
1:A:266:TYR:CE2	1:A:392:LEU:HD22	2.54	0.42
1:C:197:ARG:HB3	1:C:244:TYR:CE2	2.54	0.42
1:C:271:ALA:HB3	1:C:345:TYR:HB2	2.01	0.42
1:C:435:THR:HG23	3:C:173:HOH:O	2.18	0.42
2:B:78:LEU:HD21	2:B:117:SER:HB3	2.02	0.42
1:A:210:LYS:HE2	2:B:34:GLU:HA	2.02	0.42
1:A:400:LYS:HG2	1:A:401:ASP:N	2.35	0.42
2:D:4:VAL:HG13	2:D:60:PRO:HG3	2.01	0.42
1:A:380:TRP:CH2	1:A:382:GLY:HA2	2.55	0.42
2:D:39:ARG:HB3	2:D:52:VAL:CG2	2.44	0.42
2:D:8:LEU:HD13	2:D:11:TYR:CD1	2.54	0.42
1:A:237:VAL:HG12	1:A:242:GLU:HB3	2.02	0.41
2:B:15:PHE:CD2	2:B:15:PHE:N	2.87	0.41
2:B:117:SER:HB2	3:B:191:HOH:O	2.20	0.41
2:B:9:GLU:OE1	2:B:9:GLU:N	2.48	0.41
2:B:76:ALA:HB1	2:B:80:MSE:CE	2.51	0.41
2:B:99:GLU:OE2	2:B:115:ARG:NH1	2.54	0.41
2:D:59:TYR:HB2	2:D:60:PRO:CD	2.51	0.41
2:D:131:ILE:O	2:D:135:MSE:HG3	2.21	0.41
1:A:197:ARG:NH2	1:A:224:ALA:HB3	2.36	0.40
1:A:246:MSE:HE3	1:A:252:TYR:CB	2.50	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:350:THR:HG23	1:C:394:PHE:CE2	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:383:LYS:CG	2:D:64:LYS:O[4_545]	2.04	0.16

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	248/268 (92%)	237 (96%)	11 (4%)	0	100	100
1	C	248/268 (92%)	239 (96%)	9 (4%)	0	100	100
2	B	159/161 (99%)	152 (96%)	7 (4%)	0	100	100
2	D	158/161 (98%)	147 (93%)	11 (7%)	0	100	100
All	All	813/858 (95%)	775 (95%)	38 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	216/225 (96%)	191 (88%)	25 (12%)	4 15
1	C	216/225 (96%)	203 (94%)	13 (6%)	16 44
2	B	142/134 (106%)	131 (92%)	11 (8%)	10 31
2	D	141/134 (105%)	125 (89%)	16 (11%)	4 16
All	All	715/718 (100%)	650 (91%)	65 (9%)	7 24

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	196	GLU
1	A	197	ARG
1	A	199	LEU
1	A	200	PHE
1	A	210	LYS
1	A	212	VAL
1	A	237	VAL
1	A	256	LYS
1	A	262	LEU
1	A	264	GLN
1	A	282	LEU
1	A	309	MSE
1	A	314	THR
1	A	335	LYS
1	A	340	VAL
1	A	351	LYS
1	A	360	ARG
1	A	374	ILE
1	A	388	SER
1	A	392	LEU
1	A	403	LYS
1	A	417	PHE
1	A	428	ASN
1	A	432	LYS
1	A	444	GLN
2	B	6	LYS
2	B	18	GLU
2	B	34	GLU
2	B	64	LYS
2	B	91	GLN
2	B	103	LEU
2	B	109	LEU
2	B	113	LEU

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Mol	Chain	Res	Type
2	B	115	ARG
2	B	116	GLN
2	B	126	LEU
1	C	196	GLU
1	C	197	ARG
1	C	203	LYS
1	C	212	VAL
1	C	246	MSE
1	C	249	SER
1	C	262	LEU
1	C	282	LEU
1	C	300	LEU
1	C	333	SER
1	C	346	ASP
1	C	383	LYS
1	C	444	GLN
2	D	13	LYS
2	D	22	ASN
2	D	26	LYS
2	D	37	ILE
2	D	40	ARG
2	D	41	ASP
2	D	42	SER
2	D	54	LEU
2	D	73	ILE
2	D	80	MSE
2	D	81	SER
2	D	103	LEU
2	D	109	LEU
2	D	116	GLN
2	D	127	LYS
2	D	154	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	264	GLN
1	A	428	ASN
2	B	91	GLN
2	B	116	GLN
2	B	121	ASN
1	C	352	ASN

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Mol	Chain	Res	Type
1	C	444	GLN
2	D	22	ASN
2	D	97	ASN
2	D	116	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	245/268 (91%)	-0.39	1 (0%) 89 85	11, 21, 30, 32	0
1	C	245/268 (91%)	-0.25	4 (1%) 70 63	11, 24, 30, 35	0
2	B	153/161 (95%)	-0.46	0 100 100	9, 20, 26, 28	0
2	D	153/161 (95%)	-0.34	0 100 100	6, 19, 29, 33	0
All	All	796/858 (92%)	-0.35	5 (0%) 85 81	6, 22, 30, 35	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	235	SER	3.5
1	C	315	ASP	2.8
1	A	315	ASP	2.7
1	C	294	GLY	2.5
1	C	257	GLU	2.3

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

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

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.