



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:39 AM GMT

PDB ID : 2XVU
Title : HUMAN SERUM ALBUMIN COMPLEXED WITH DANSYL-L-ASPARAGINE
Authors : Zunszain, P.A.; Curry, S.
Deposited on : 2010-10-28
Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

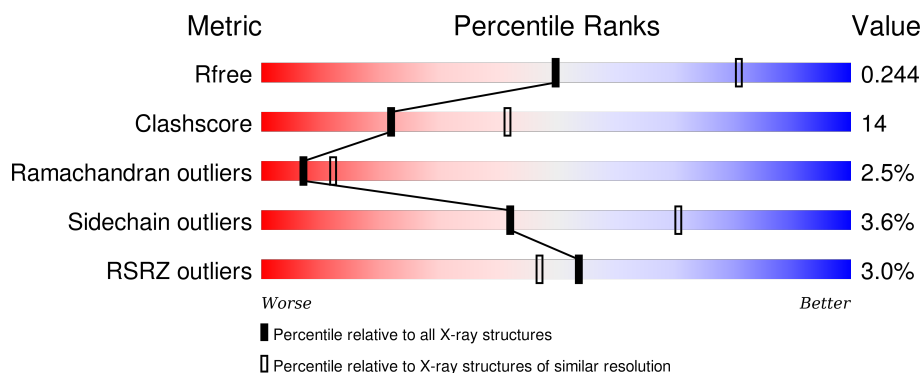
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	585	<div> <div>3%</div> <div>69%</div> <div>24%</div> <div>• •</div> </div>
1	B	585	<div> <div>3%</div> <div>67%</div> <div>26%</div> <div>• 5%</div> </div>

2 Entry composition [i](#)

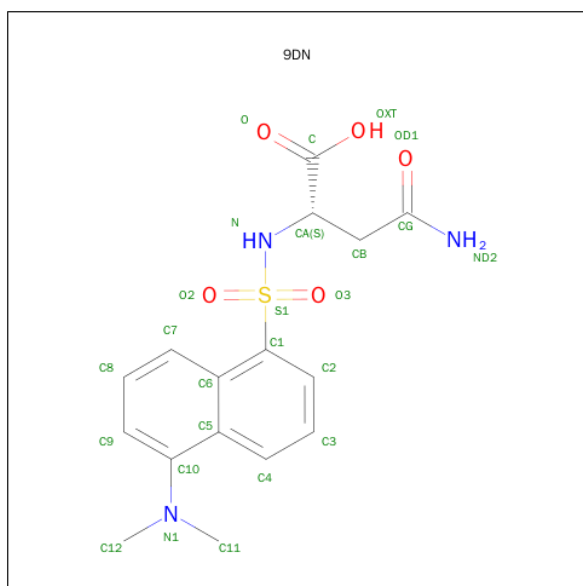
There are 3 unique types of molecules in this entry. The entry contains 8366 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SERUM ALBUMIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	560	Total	C	N	O	S	0	0	0
			4148	2626	689	794	39			
1	B	557	Total	C	N	O	S	0	0	0
			4083	2578	682	783	40			

- Molecule 2 is DANSYL-L-ASPARAGINE (three-letter code: 9DN) (formula: $C_{16}H_{19}N_3O_5S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			25	16	3	5	1		
2	A	1	Total	C	N	O	S	0	0
			25	16	3	5	1		
2	B	1	Total	C	N	O	S	0	0
			25	16	3	5	1		
2	B	1	Total	C	N	O	S	0	0
			25	16	3	5	1		

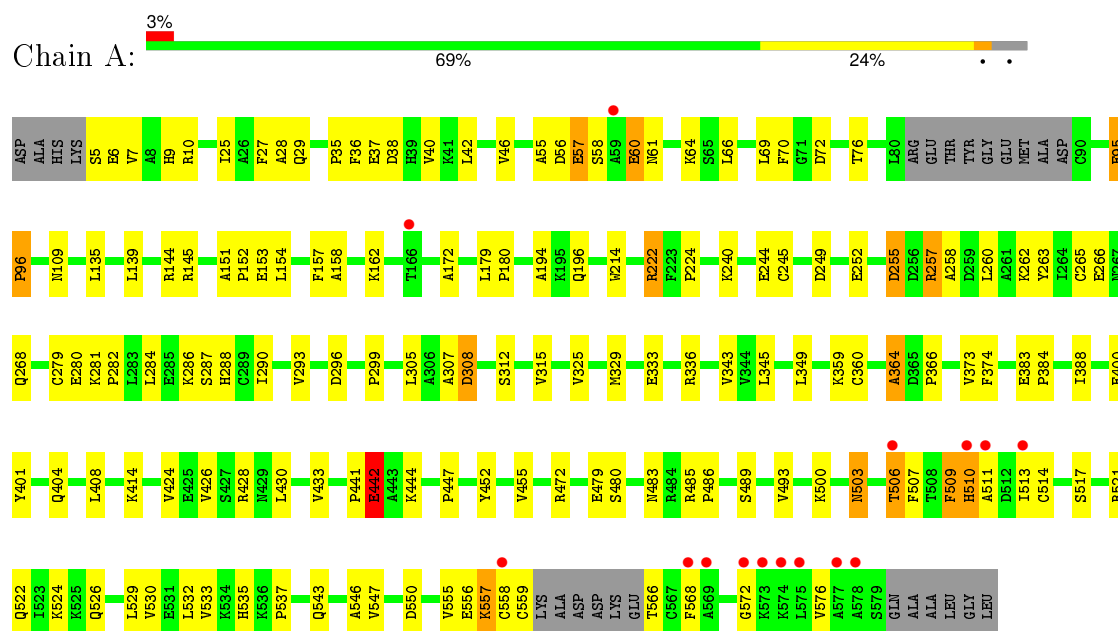
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	17	Total 17	O 17	0	0
3	B	18	Total 18	O 18	0	0

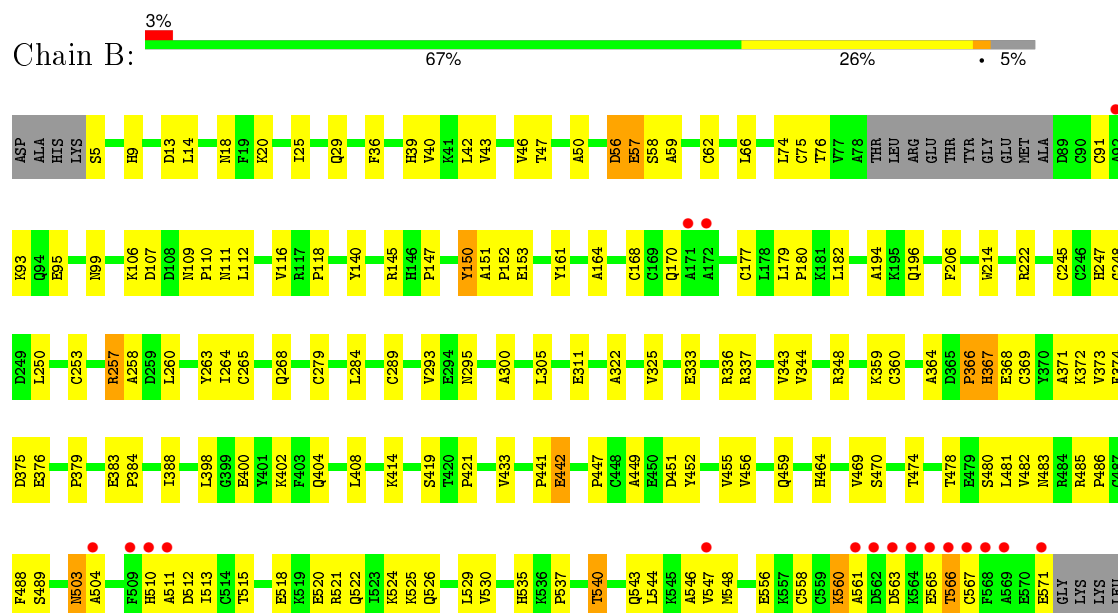
3 Residue-property plots

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: SERUM ALBUMIN



• Molecule 1: SERUM ALBUMIN



VAL
ALA
ALA
SER
GLN
ALA
ALA
LEU
GLY
LEU

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	54.85Å 55.25Å 119.73Å 81.34° 90.93° 64.60°	Depositor
Resolution (Å)	35.12 – 2.60 35.12 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.7 (35.12-2.60) 91.6 (35.12-2.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 2.61Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.205 , 0.244 0.205 , 0.244	Depositor DCC
R_{free} test set	1791 reflections (4.77%)	DCC
Wilson B-factor (Å ²)	55.6	Xtriage
Anisotropy	0.314	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 61.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 38418 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8366	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.69% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 9DN

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	0/4229	0.60	0/5759
1	B	0.36	0/4162	0.58	0/5674
All	All	0.37	0/8391	0.59	0/11433

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	4148	0	3750	103	0
1	B	4083	0	3660	117	0
2	A	50	0	36	1	0
2	B	50	0	36	1	0
3	A	17	0	0	1	0
3	B	18	0	0	2	0
All	All	8366	0	7482	220	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 14.

All (220) 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:540:THR:HG23	1:B:543:GLN:HB3	1.50	0.94
1:B:483:ASN:O	1:B:486:PRO:HD2	1.77	0.84
1:A:383:GLU:HB3	1:A:384:PRO:HD3	1.57	0.84
1:A:265:CYS:O	1:A:268:GLN:HG3	1.79	0.83
1:B:433:VAL:HG22	1:B:452:TYR:CD2	2.12	0.83
1:B:151:ALA:HB3	1:B:152:PRO:HD3	1.62	0.81
1:A:511:ALA:HA	1:A:514:CYS:SG	2.20	0.81
1:B:414:LYS:HE3	1:B:488:PHE:O	1.82	0.78
1:A:151:ALA:HB3	1:A:152:PRO:HD3	1.65	0.78
1:B:265:CYS:O	1:B:268:GLN:HG3	1.82	0.77
1:A:66:LEU:O	1:A:70:PHE:HD1	1.68	0.77
1:B:419:SER:OG	1:B:421:PRO:HD2	1.84	0.76
1:B:567:CYS:SG	1:B:571:GLU:HB2	2.26	0.76
1:B:59:ALA:HB3	1:B:62:CYS:SG	2.25	0.76
1:B:383:GLU:HB3	1:B:384:PRO:HD3	1.68	0.75
1:A:433:VAL:HG22	1:A:452:TYR:CD2	2.23	0.73
1:A:364:ALA:O	1:A:366:PRO:HD3	1.88	0.73
1:B:540:THR:HG23	1:B:543:GLN:CB	2.19	0.73
1:A:510:HIS:CB	1:A:568:PHE:HA	2.20	0.72
1:A:424:VAL:O	1:A:428:ARG:HG3	1.90	0.71
1:B:95:GLU:OE1	1:B:99:ASN:HB2	1.90	0.70
1:A:145:ARG:HG3	1:A:145:ARG:HH11	1.56	0.70
1:A:384:PRO:O	1:A:388:ILE:HG12	1.93	0.68
1:A:509:PHE:O	1:A:510:HIS:CB	2.41	0.67
1:A:224:PRO:HD2	1:A:296:ASP:HB3	1.77	0.67
1:A:61:ASN:HD22	1:A:64:LYS:HE3	1.60	0.67
1:B:106:LYS:HD3	1:B:147:PRO:HB2	1.78	0.65
1:B:116:VAL:O	1:B:118:PRO:HD3	1.97	0.65
1:A:483:ASN:O	1:A:486:PRO:HD2	1.97	0.65
1:B:305:LEU:HD21	1:B:333:GLU:HB3	1.79	0.64
1:A:95:GLU:CB	1:A:96:PRO:HD3	2.28	0.63
1:A:36:PHE:O	1:A:40:VAL:HG23	1.99	0.63
1:A:483:ASN:C	1:A:486:PRO:HD2	2.18	0.63
1:A:359:LYS:HG3	1:A:360:CYS:N	2.13	0.63
1:B:348:ARG:HG3	1:B:482:VAL:HG12	1.81	0.63
1:A:139:LEU:HD21	1:A:158:ALA:HB2	1.80	0.62
1:A:373:VAL:HG13	1:A:374:PHE:HD1	1.65	0.61
1:A:566:THR:C	1:A:568:PHE:N	2.52	0.61
1:B:206:PHE:CD2	1:B:481:LEU:HD22	2.36	0.61
1:B:179:LEU:HB2	1:B:180:PRO:HD3	1.84	0.60
1:B:483:ASN:C	1:B:486:PRO:HD2	2.20	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:260:LEU:HD23	2:A:2001:9DN:H122	1.82	0.60
1:B:398:LEU:HB3	1:B:402:LYS:HB2	1.83	0.60
1:A:566:THR:C	1:A:568:PHE:H	2.04	0.58
1:B:260:LEU:O	1:B:264:ILE:HG13	2.02	0.58
1:B:348:ARG:HG3	1:B:482:VAL:CG1	2.33	0.58
1:B:247:HIS:O	1:B:247:HIS:ND1	2.37	0.58
1:A:305:LEU:HD21	1:A:333:GLU:HB3	1.84	0.58
1:B:39:HIS:O	1:B:43:VAL:HG23	2.03	0.58
1:B:25:ILE:O	1:B:29:GLN:HG3	2.04	0.57
1:A:179:LEU:HB2	1:A:180:PRO:HD3	1.85	0.57
1:B:206:PHE:CE2	1:B:481:LEU:HD13	2.40	0.57
1:B:373:VAL:HG13	1:B:374:PHE:HD1	1.68	0.57
1:A:547:VAL:HA	1:A:550:ASP:OD2	2.04	0.57
1:B:529:LEU:HD22	1:B:548:MET:HE2	1.87	0.57
1:A:400:GLU:O	1:A:404:GLN:HG3	2.05	0.57
1:A:25:ILE:HD13	1:A:154:LEU:HD23	1.86	0.56
1:B:518:GLU:O	1:B:522:GLN:HG3	2.05	0.56
1:B:145:ARG:HH11	1:B:145:ARG:HG3	1.71	0.55
1:B:257:ARG:HD3	1:B:258:ALA:N	2.21	0.55
1:B:544:LEU:O	1:B:547:VAL:HG12	2.07	0.55
1:B:464:HIS:CE1	1:B:469:VAL:H	2.24	0.55
1:B:5:SER:HA	1:B:62:CYS:O	2.06	0.55
1:B:214:TRP:CD1	1:B:343:VAL:HG11	2.41	0.55
1:A:66:LEU:O	1:A:70:PHE:CD1	2.57	0.54
1:A:257:ARG:HD3	1:A:258:ALA:N	2.22	0.54
1:A:61:ASN:HB3	1:A:64:LYS:HD2	1.90	0.54
1:A:500:LYS:O	1:A:535:HIS:ND1	2.41	0.54
1:B:257:ARG:NH2	1:B:284:LEU:HD23	2.23	0.54
1:B:563:ASP:HB2	1:B:566:THR:HB	1.90	0.53
1:A:56:ASP:C	1:A:58:SER:H	2.11	0.53
1:B:400:GLU:O	1:B:404:GLN:HG3	2.08	0.53
1:B:388:ILE:HA	2:B:2002:9DN:H9	1.91	0.53
1:A:279:CYS:HA	1:A:286:LYS:HD2	1.91	0.53
1:A:249:ASP:HB3	1:A:252:GLU:CG	2.40	0.52
1:B:408:LEU:HD11	1:B:526:GLN:HB3	1.91	0.52
1:A:38:ASP:O	1:A:42:LEU:HG	2.10	0.52
1:B:337:ARG:HG3	1:B:337:ARG:HH11	1.75	0.52
1:A:281:LYS:HB2	1:A:282:PRO:CD	2.40	0.52
1:B:257:ARG:HH22	1:B:284:LEU:HD23	1.74	0.52
1:A:42:LEU:O	1:A:46:VAL:HG23	2.10	0.52
1:A:61:ASN:HD22	1:A:64:LYS:CE	2.21	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:LEU:O	1:B:46:VAL:HG23	2.09	0.51
1:A:555:VAL:O	1:A:559:CYS:HB2	2.10	0.51
1:B:376:GLU:O	1:B:379:PRO:HD2	2.10	0.51
1:B:56:ASP:O	1:B:57:GLU:C	2.49	0.51
1:A:145:ARG:NH1	1:A:145:ARG:HG3	2.23	0.51
1:B:39:HIS:HD2	1:B:140:TYR:HE1	1.59	0.51
1:A:290:ILE:O	1:A:293:VAL:HG12	2.11	0.51
1:B:150:TYR:HD2	1:B:153:GLU:HB2	1.76	0.51
1:B:372:LYS:N	3:B:3009:HOH:O	2.25	0.50
1:A:262:LYS:O	1:A:266:GLU:HG3	2.11	0.50
1:A:572:GLY:O	1:A:576:VAL:HG23	2.11	0.50
1:B:483:ASN:O	1:B:486:PRO:CD	2.56	0.50
1:B:520:GLU:O	1:B:524:LYS:HG2	2.11	0.50
1:B:20:LYS:HE3	1:B:47:THR:HG21	1.94	0.50
1:B:561:ALA:HB3	1:B:563:ASP:OD1	2.11	0.50
1:B:543:GLN:O	1:B:546:ALA:HB3	2.12	0.50
1:B:485:ARG:HB3	1:B:486:PRO:HD3	1.92	0.50
1:A:373:VAL:HG13	1:A:374:PHE:CD1	2.45	0.50
1:B:289:CYS:O	1:B:293:VAL:HG13	2.11	0.50
1:B:398:LEU:HB3	1:B:402:LYS:CB	2.42	0.49
1:A:522:GLN:O	1:A:526:GLN:HG3	2.13	0.49
1:B:367:HIS:O	1:B:371:ALA:HB2	2.12	0.49
1:A:506:THR:HG22	1:A:507:PHE:CD1	2.48	0.49
1:B:511:ALA:C	1:B:513:ILE:H	2.14	0.49
1:A:485:ARG:HB3	1:A:486:PRO:HD3	1.94	0.49
1:A:345:LEU:O	1:A:349:LEU:HG	2.12	0.49
1:B:529:LEU:HD22	1:B:548:MET:CE	2.42	0.49
1:A:513:ILE:HG13	1:A:514:CYS:N	2.26	0.49
1:B:47:THR:O	1:B:50:ALA:HB3	2.12	0.48
1:A:414:LYS:O	1:A:472:ARG:NH1	2.46	0.48
1:A:25:ILE:CD1	1:A:154:LEU:HD23	2.44	0.48
1:B:447:PRO:O	1:B:451:ASP:HB2	2.13	0.48
1:B:194:ALA:HB1	1:B:455:VAL:CG1	2.43	0.48
1:A:72:ASP:O	1:A:76:THR:HG23	2.13	0.48
1:A:521:ARG:O	1:A:524:LYS:HB2	2.14	0.48
1:B:110:PRO:HB2	1:B:112:LEU:HG	1.96	0.48
1:B:111:ASN:O	1:B:112:LEU:HD23	2.14	0.48
1:B:511:ALA:C	1:B:513:ILE:N	2.67	0.48
1:A:7:VAL:HG21	1:A:69:LEU:CD1	2.44	0.48
1:B:364:ALA:O	1:B:366:PRO:HD3	2.14	0.48
1:A:35:PRO:HD2	1:A:38:ASP:OD2	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:36:PHE:O	1:B:40:VAL:HG23	2.14	0.47
1:B:414:LYS:CE	1:B:488:PHE:O	2.58	0.47
1:B:9:HIS:HD2	1:B:13:ASP:OD1	1.98	0.47
1:A:66:LEU:HD12	1:A:252:GLU:OE2	2.15	0.47
1:B:408:LEU:HD22	1:B:530:VAL:CG2	2.44	0.47
1:A:153:GLU:OE2	1:A:288:HIS:ND1	2.41	0.47
1:B:322:ALA:HB1	1:B:325:VAL:HB	1.96	0.47
1:B:161:TYR:O	1:B:164:ALA:HB3	2.14	0.47
1:B:474:THR:O	1:B:478:THR:HG23	2.15	0.47
1:A:9:HIS:C	1:A:9:HIS:ND1	2.68	0.47
1:A:308:ASP:N	1:A:308:ASP:OD1	2.48	0.47
1:A:543:GLN:O	1:A:546:ALA:HB3	2.15	0.47
1:B:535:HIS:C	1:B:537:PRO:HD3	2.34	0.47
1:B:384:PRO:O	1:B:388:ILE:HG12	2.15	0.46
1:A:359:LYS:CG	1:A:360:CYS:N	2.78	0.46
1:A:281:LYS:HB2	1:A:282:PRO:HD2	1.97	0.46
1:B:39:HIS:HD2	1:B:140:TYR:CE1	2.33	0.46
1:A:257:ARG:HD3	1:A:257:ARG:C	2.36	0.46
1:A:408:LEU:HD22	1:A:530:VAL:CG2	2.46	0.46
1:A:401:TYR:HE1	1:A:522:GLN:HB3	1.80	0.46
1:B:222:ARG:HA	1:B:295:ASN:OD1	2.15	0.46
1:A:25:ILE:O	1:A:29:GLN:HG3	2.15	0.46
1:B:74:LEU:C	1:B:76:THR:H	2.20	0.46
1:A:325:VAL:O	1:A:329:MET:HG3	2.15	0.46
1:B:456:VAL:O	1:B:459:GLN:HB2	2.16	0.45
1:B:305:LEU:CD2	1:B:333:GLU:HB3	2.45	0.45
1:A:222:ARG:O	1:A:224:PRO:HD3	2.16	0.45
1:A:408:LEU:HD11	1:A:526:GLN:HB3	1.97	0.45
1:A:222:ARG:C	1:A:224:PRO:HD3	2.37	0.45
1:B:91:CYS:C	1:B:93:LYS:H	2.18	0.45
1:B:388:ILE:HD12	1:B:449:ALA:CB	2.46	0.45
1:A:5:SER:HB3	1:A:57:GLU:HG3	1.99	0.45
1:A:535:HIS:C	1:A:537:PRO:HD3	2.38	0.45
1:A:257:ARG:HH22	1:A:284:LEU:HD23	1.81	0.45
1:B:91:CYS:C	1:B:93:LYS:N	2.69	0.44
1:B:510:HIS:CD2	1:B:512:ASP:H	2.35	0.44
1:B:253:CYS:O	1:B:257:ARG:HB3	2.17	0.44
1:A:240:LYS:O	1:A:244:GLU:HG3	2.17	0.44
1:B:66:LEU:HG	3:B:3003:HOH:O	2.17	0.44
1:A:503:ASN:O	1:A:506:THR:HB	2.17	0.44
1:B:91:CYS:O	1:B:93:LYS:N	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:464:HIS:CE1	1:B:470:SER:H	2.36	0.44
1:B:263:TYR:C	1:B:263:TYR:CD1	2.91	0.44
1:B:196:GLN:HA	1:B:196:GLN:NE2	2.33	0.44
1:B:503:ASN:HA	1:B:503:ASN:HD22	1.54	0.43
1:B:511:ALA:O	1:B:513:ILE:N	2.51	0.43
1:A:153:GLU:O	1:A:157:PHE:HD2	2.01	0.43
1:A:214:TRP:CD1	1:A:343:VAL:HG11	2.53	0.43
1:B:359:LYS:CG	1:B:360:CYS:N	2.81	0.43
1:B:336:ARG:HD3	1:B:337:ARG:HH12	1.83	0.43
1:B:558:CYS:C	1:B:560:LYS:H	2.21	0.43
1:A:566:THR:O	1:A:568:PHE:N	2.51	0.43
1:B:560:LYS:O	1:B:560:LYS:HG3	2.18	0.43
1:A:10:ARG:NH1	1:A:255:ASP:OD2	2.44	0.43
1:B:14:LEU:O	1:B:18:ASN:HB2	2.19	0.42
1:A:444:LYS:O	1:A:447:PRO:HD2	2.20	0.42
1:B:540:THR:CG2	1:B:543:GLN:HB3	2.34	0.42
1:A:28:ALA:O	1:A:144:ARG:HA	2.19	0.42
1:A:426:VAL:O	1:A:430:LEU:HG	2.19	0.42
1:B:168:CYS:SG	1:B:177:CYS:C	2.98	0.42
1:A:27:PHE:HE1	1:A:70:PHE:HD2	1.68	0.42
1:B:145:ARG:HG3	1:B:145:ARG:NH1	2.35	0.42
1:B:366:PRO:O	1:B:368:GLU:N	2.52	0.42
1:B:366:PRO:O	1:B:369:CYS:N	2.52	0.42
1:B:388:ILE:HD12	1:B:449:ALA:HB3	2.01	0.42
1:B:257:ARG:HD3	1:B:257:ARG:C	2.40	0.42
1:A:135:LEU:HD11	1:A:162:LYS:CD	2.49	0.41
1:B:248:GLY:O	1:B:250:LEU:N	2.52	0.41
1:B:107:ASP:HB3	1:B:110:PRO:HG3	2.02	0.41
1:B:109:ASN:N	1:B:110:PRO:HD3	2.34	0.41
1:B:168:CYS:C	1:B:170:GLN:H	2.23	0.41
1:B:521:ARG:O	1:B:524:LYS:HB2	2.21	0.41
1:A:95:GLU:CB	1:A:96:PRO:CD	2.98	0.41
1:A:257:ARG:HH22	1:A:284:LEU:CD2	2.32	0.41
1:A:6:GLU:O	1:A:9:HIS:HB3	2.21	0.41
1:A:60:GLU:OE1	1:A:60:GLU:HA	2.21	0.41
1:A:529:LEU:O	1:A:533:VAL:HG23	2.20	0.41
1:A:249:ASP:HB3	1:A:252:GLU:CD	2.40	0.41
1:A:27:PHE:CE1	1:A:70:PHE:HD2	2.38	0.41
1:A:281:LYS:HE2	1:A:281:LYS:HB3	1.82	0.41
1:A:506:THR:HG22	1:A:507:PHE:CE1	2.56	0.41
1:B:57:GLU:O	1:B:59:ALA:N	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:179:LEU:CD2	1:B:179:LEU:H	2.34	0.41
1:A:556:GLU:HG3	1:A:557:LYS:H	1.86	0.41
1:A:194:ALA:HB1	1:A:455:VAL:CG1	2.51	0.41
1:A:441:PRO:O	1:A:442:GLU:C	2.59	0.41
1:B:556:GLU:O	1:B:556:GLU:HG2	2.21	0.40
1:A:224:PRO:HB2	1:A:299:PRO:HD3	2.03	0.40
1:B:260:LEU:O	1:B:260:LEU:HD23	2.21	0.40
1:B:522:GLN:HA	1:B:525:LYS:HD3	2.03	0.40
1:B:544:LEU:C	1:B:546:ALA:N	2.74	0.40
1:B:305:LEU:HA	1:B:305:LEU:HD23	1.92	0.40
1:B:441:PRO:O	1:B:442:GLU:C	2.60	0.40
1:A:196:GLN:NE2	3:A:3007:HOH:O	2.54	0.40
1:A:333:GLU:HA	1:A:336:ARG:HD2	2.03	0.40
1:A:557:LYS:HG2	1:A:558:CYS:N	2.37	0.40
1:A:263:TYR:CD1	1:A:263:TYR:C	2.94	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	554/585 (95%)	496 (90%)	43 (8%)	15 (3%)	6	10
1	B	553/585 (94%)	481 (87%)	59 (11%)	13 (2%)	7	13
All	All	1107/1170 (95%)	977 (88%)	102 (9%)	28 (2%)	7	12

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	95	GLU
1	A	510	HIS
1	B	57	GLU

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Mol	Chain	Res	Type
1	B	58	SER
1	B	565	GLU
1	A	55	ALA
1	A	172	ALA
1	A	509	PHE
1	A	517	SER
1	B	279	CYS
1	B	300	ALA
1	B	367	HIS
1	B	566	THR
1	A	364	ALA
1	B	75	CYS
1	B	150	TYR
1	B	504	ALA
1	B	560	LYS
1	A	57	GLU
1	A	60	GLU
1	A	442	GLU
1	A	312	SER
1	A	315	VAL
1	B	442	GLU
1	A	307	ALA
1	A	557	LYS
1	A	96	PRO
1	B	366	PRO

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	403/511 (79%)	386 (96%)	17 (4%)	36	65
1	B	392/511 (77%)	380 (97%)	12 (3%)	47	76
All	All	795/1022 (78%)	766 (96%)	29 (4%)	42	71

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	GLU
1	A	109	ASN
1	A	222	ARG
1	A	245	CYS
1	A	255	ASP
1	A	257	ARG
1	A	280	GLU
1	A	287	SER
1	A	308	ASP
1	A	442	GLU
1	A	479	GLU
1	A	480	SER
1	A	489	SER
1	A	493	VAL
1	A	503	ASN
1	A	506	THR
1	A	532	LEU
1	B	56	ASP
1	B	182	LEU
1	B	245	CYS
1	B	257	ARG
1	B	311	GLU
1	B	344	VAL
1	B	375	ASP
1	B	480	SER
1	B	489	SER
1	B	503	ASN
1	B	515	THR
1	B	540	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	ASN
1	A	109	ASN
1	A	196	GLN
1	A	247	HIS
1	A	464	HIS
1	A	483	ASN
1	A	503	ASN
1	A	543	GLN
1	B	9	HIS
1	B	39	HIS

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Mol	Chain	Res	Type
1	B	196	GLN
1	B	464	HIS
1	B	483	ASN
1	B	503	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](#)

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	9DN	A	2001	-	23,26,26	2.21	7 (30%)	32,38,38	1.11	3 (9%)
2	9DN	A	2002	-	23,26,26	1.91	5 (21%)	32,38,38	1.23	3 (9%)
2	9DN	B	2001	-	23,26,26	2.27	8 (34%)	32,38,38	1.12	2 (6%)
2	9DN	B	2002	-	23,26,26	1.87	5 (21%)	32,38,38	1.26	5 (15%)

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	9DN	A	2001	-	-	0/19/23/23	0/2/2/2
2	9DN	A	2002	-	-	0/19/23/23	0/2/2/2
2	9DN	B	2001	-	-	0/19/23/23	0/2/2/2
2	9DN	B	2002	-	-	0/19/23/23	0/2/2/2

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2001	9DN	C10-C5	2.01	1.47	1.42
2	B	2001	9DN	C3-C2	2.02	1.43	1.38
2	A	2001	9DN	C3-C2	2.05	1.43	1.38
2	A	2001	9DN	S1-N	2.26	1.65	1.61
2	B	2001	9DN	S1-N	2.33	1.65	1.61
2	A	2002	9DN	C6-C5	2.85	1.48	1.43
2	A	2002	9DN	C9-C10	2.89	1.44	1.38
2	B	2002	9DN	C6-C5	2.97	1.48	1.43
2	A	2001	9DN	C6-C5	3.00	1.49	1.43
2	B	2002	9DN	C9-C10	3.01	1.44	1.38
2	A	2001	9DN	C9-C10	3.06	1.45	1.38
2	B	2001	9DN	C6-C5	3.21	1.49	1.43
2	B	2001	9DN	C9-C10	3.25	1.45	1.38
2	A	2002	9DN	C8-C9	3.28	1.45	1.38
2	A	2001	9DN	C8-C9	3.39	1.45	1.38
2	B	2001	9DN	C8-C9	3.61	1.46	1.38
2	B	2002	9DN	C1-C6	3.67	1.49	1.43
2	B	2002	9DN	C8-C9	3.70	1.46	1.38
2	A	2002	9DN	C1-C6	3.83	1.49	1.43
2	B	2002	9DN	C1-S1	3.87	1.82	1.77
2	A	2001	9DN	C1-C6	4.05	1.49	1.43
2	B	2001	9DN	C1-C6	4.09	1.49	1.43
2	A	2002	9DN	C1-S1	4.54	1.82	1.77
2	A	2001	9DN	C1-S1	5.94	1.84	1.77
2	B	2001	9DN	C1-S1	5.94	1.84	1.77

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2001	9DN	C12-N1-C11	-3.01	106.03	115.96
2	B	2001	9DN	C12-N1-C11	-2.98	106.10	115.96
2	B	2002	9DN	C12-N1-C11	-2.97	106.15	115.96
2	A	2002	9DN	C12-N1-C11	-2.74	106.89	115.96
2	A	2002	9DN	O3-S1-O2	-2.32	116.46	119.54
2	B	2002	9DN	O3-S1-O2	-2.24	116.57	119.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2001	9DN	O3-S1-O2	-2.15	116.69	119.54
2	B	2002	9DN	C4-C5-C10	-2.06	119.23	122.49
2	A	2001	9DN	C10-C5-C6	2.06	121.64	119.35
2	B	2002	9DN	O2-S1-C1	2.07	111.53	108.01
2	B	2001	9DN	C10-C5-C6	2.14	121.72	119.35
2	A	2002	9DN	C10-C5-C6	2.19	121.78	119.35
2	B	2002	9DN	C10-C5-C6	2.36	121.97	119.35

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2001	9DN	1	0
2	B	2002	9DN	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	560/585 (95%)	-0.28	15 (2%) 58 51	29, 59, 121, 155	0
1	B	557/585 (95%)	-0.25	18 (3%) 51 44	26, 62, 113, 146	0
All	All	1117/1170 (95%)	-0.26	33 (2%) 54 47	26, 61, 118, 155	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	562	ASP	6.1
1	A	569	ALA	6.1
1	B	564	LYS	4.9
1	A	510	HIS	4.6
1	A	511	ALA	4.1
1	B	566	THR	3.9
1	A	568	PHE	3.6
1	B	563	ASP	3.4
1	B	561	ALA	3.4
1	A	578	ALA	3.2
1	B	171	ALA	3.1
1	A	513	ILE	3.0
1	B	510	HIS	3.0
1	B	569	ALA	3.0
1	B	565	GLU	2.8
1	A	574	LYS	2.8
1	B	571	GLU	2.7
1	B	172	ALA	2.5
1	B	567	CYS	2.5
1	A	573	LYS	2.5
1	A	572	GLY	2.5
1	A	575	LEU	2.5
1	A	558	CYS	2.5
1	A	577	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	568	PHE	2.4
1	B	511	ALA	2.4
1	A	166	THR	2.3
1	A	506	THR	2.2
1	B	547	VAL	2.1
1	A	59	ALA	2.1
1	B	509	PHE	2.1
1	B	92	ALA	2.1
1	B	504	ALA	2.1

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	9DN	A	2002	25/25	0.96	0.19	1.41	52,56,71,74	0
2	9DN	B	2002	25/25	0.95	0.20	1.13	60,63,70,71	0
2	9DN	A	2001	25/25	0.94	0.18	0.77	68,72,74,74	0
2	9DN	B	2001	25/25	0.93	0.19	0.73	69,76,81,83	0

6.5 Other polymers [i](#)

There are no such residues in this entry.