



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 01:37 PM GMT

PDB ID : 2BXF
Title : HUMAN SERUM ALBUMIN COMPLEXED WITH DIAZEPAM
Authors : Ghuman, J.; Zunszain, P.A.; Petitpas, I.; Bhattacharya, A.A.; Curry, S.
Deposited on : 2005-07-26
Resolution : 2.95 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

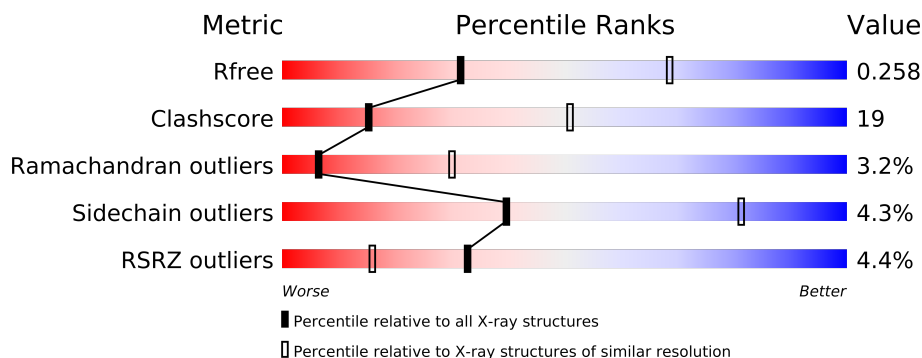
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.95 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}	66092	1587 (3.00-2.92)
Clashscore	79885	2029 (3.00-2.92)
Ramachandran outliers	78287	1955 (3.00-2.92)
Sidechain outliers	78261	1958 (3.00-2.92)
RSRZ outliers	66119	1588 (3.00-2.92)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	585	
1	B	585	

2 Entry composition i

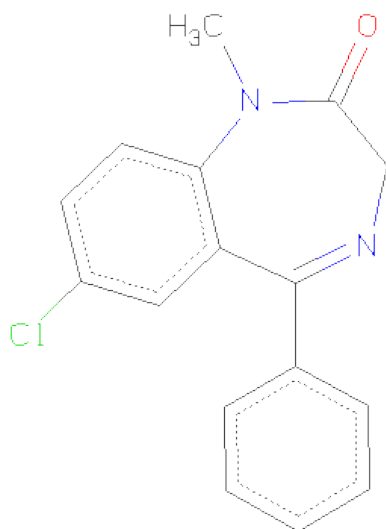
There are 2 unique types of molecules in this entry. The entry contains 8615 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

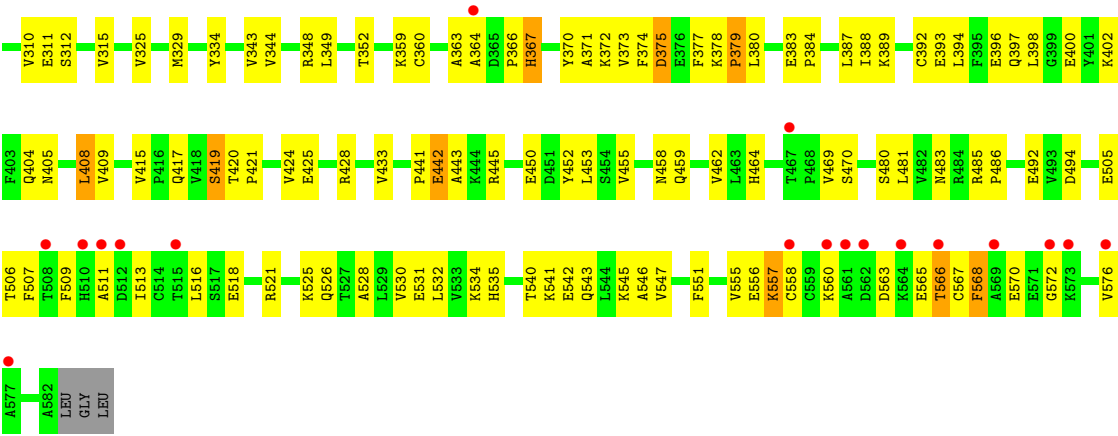
- Molecule 1 is a protein called SERUM ALBUMIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	578	Total	C	N	O	S	0	0	0
			4282	2700	713	828	41			
1	B	578	Total	C	N	O	S	0	0	0
			4293	2711	714	828	40			

- Molecule 2 is 7-CHLORO-1-METHYL-5-PHENYL-1,3-DIHYDRO-2H-1,4-BENZODIAZEPIN-2-ONE (three-letter code: DZP) (formula: C₁₆H₁₃ClN₂O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	0	0
			20	16	1	2	1		
2	B	1	Total	C	Cl	N	O	0	0
			20	16	1	2	1		



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.29Å 55.78Å 120.51Å 81.17° 90.22° 65.93°	Depositor
Resolution (Å)	22.59 – 2.95 22.59 – 2.95	Depositor EDS
% Data completeness (in resolution range)	96.7 (22.59-2.95) 96.7 (22.59-2.95)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.62 (at 2.94Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.215 , 0.262 0.213 , 0.258	Depositor DCC
R_{free} test set	1243 reflections (4.96%)	DCC
Wilson B-factor (Å ²)	71.2	Xtriage
Anisotropy	0.308	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 65.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 27216 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8615	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0/4366	0.61	0/5949
1	B	0.38	0/4377	0.60	0/5961
All	All	0.38	0/8743	0.60	0/11910

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4282	0	3874	153	0
1	B	4293	0	3908	156	0
2	A	20	0	13	0	0
2	B	20	0	13	1	0
All	All	8615	0	7808	309	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 19.

All (309) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:540:THR:HG22	1:B:541:LYS:H	1.03	1.08
1:B:383:GLU:HB3	1:B:384:PRO:HD3	1.48	0.94
1:A:540:THR:HG22	1:A:541:LYS:H	1.33	0.94
1:B:424:VAL:O	1:B:428:ARG:HG3	1.67	0.93
1:A:424:VAL:O	1:A:428:ARG:HG3	1.70	0.92
1:B:540:THR:HG22	1:B:541:LYS:N	1.87	0.87
1:B:394:LEU:O	1:B:397:GLN:HG2	1.78	0.83
1:A:214:TRP:CD1	1:A:343:VAL:HG11	2.13	0.82
1:B:540:THR:CG2	1:B:541:LYS:H	1.85	0.82
1:B:433:VAL:HG22	1:B:452:TYR:HD2	1.44	0.82
1:A:151:ALA:HB3	1:A:152:PRO:HD3	1.61	0.82
1:B:419:SER:OG	1:B:421:PRO:HD2	1.81	0.80
1:B:206:PHE:CE2	1:B:481:LEU:HD13	2.19	0.78
1:A:23:VAL:HG13	1:A:70:PHE:HE2	1.49	0.77
1:B:480:SER:OG	1:B:483:ASN:HB2	1.84	0.77
1:B:483:ASN:O	1:B:486:PRO:HD2	1.84	0.76
1:A:152:PRO:HB2	1:A:257:ARG:HH11	1.49	0.76
1:B:151:ALA:HB3	1:B:152:PRO:HD3	1.68	0.76
1:A:224:PRO:HB2	1:A:299:PRO:HD3	1.66	0.75
1:B:135:LEU:HD11	1:B:162:LYS:HD3	1.68	0.75
1:B:16:GLU:O	1:B:20:LYS:HG2	1.86	0.75
1:A:483:ASN:O	1:A:486:PRO:HD2	1.86	0.74
1:B:383:GLU:OE2	1:B:387:LEU:HD11	1.88	0.74
1:A:540:THR:HG22	1:A:541:LYS:N	2.04	0.72
1:A:120:VAL:HG21	1:A:175:ALA:HA	1.72	0.70
1:B:531:GLU:O	1:B:535:HIS:HD2	1.74	0.70
1:A:420:THR:O	1:A:424:VAL:HG23	1.91	0.69
1:B:120:VAL:HG21	1:B:175:ALA:HA	1.74	0.69
1:B:325:VAL:HG12	1:B:329:MET:CE	2.24	0.68
1:B:194:ALA:HB1	1:B:455:VAL:CG1	2.25	0.67
1:A:531:GLU:HA	1:A:531:GLU:OE1	1.93	0.67
1:B:433:VAL:HG22	1:B:452:TYR:CD2	2.28	0.66
1:B:378:LYS:HB2	1:B:379:PRO:HD3	1.77	0.66
1:A:183:ASP:O	1:A:186:ARG:HG2	1.96	0.66
1:B:392:CYS:O	1:B:396:GLU:HG3	1.95	0.66
1:A:61:ASN:HB3	1:A:64:LYS:HD3	1.78	0.65
1:A:168:CYS:SG	1:A:177:CYS:C	2.75	0.65
1:B:30:TYR:HE1	1:B:103:LEU:HD23	1.61	0.65
1:A:464:HIS:HE1	1:A:470:SER:H	1.43	0.65
1:A:152:PRO:CB	1:A:257:ARG:HH11	2.09	0.65
1:B:39:HIS:O	1:B:43:VAL:HG23	1.97	0.64
1:A:512:ASP:O	1:A:515:THR:HG22	1.97	0.64
1:A:376:GLU:O	1:A:379:PRO:HD2	1.97	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:34:CYS:HB3	1:B:39:HIS:NE2	2.14	0.63
1:B:106:LYS:HD3	1:B:147:PRO:HB2	1.79	0.63
1:B:359:LYS:HG3	1:B:360:CYS:N	2.13	0.63
1:B:59:ALA:HB3	1:B:62:CYS:SG	2.39	0.62
1:A:117:ARG:HH12	1:A:182:LEU:HB3	1.63	0.62
1:A:567:CYS:O	1:A:571:GLU:HB2	1.99	0.62
1:B:521:ARG:NH1	1:B:521:ARG:HB2	2.15	0.62
1:B:325:VAL:HG12	1:B:329:MET:HE2	1.81	0.62
1:A:194:ALA:HB1	1:A:455:VAL:CG1	2.30	0.62
1:A:539:ALA:HB1	1:A:543:GLN:HG3	1.81	0.62
1:B:464:HIS:CE1	1:B:469:VAL:H	2.17	0.62
1:B:572:GLY:O	1:B:576:VAL:HG23	2.00	0.61
1:A:540:THR:CG2	1:A:541:LYS:H	2.12	0.61
1:A:152:PRO:HB2	1:A:257:ARG:NH1	2.14	0.61
1:A:206:PHE:CD2	1:A:481:LEU:HD22	2.35	0.61
1:A:117:ARG:NH1	1:A:182:LEU:HB3	2.15	0.61
1:B:516:LEU:O	1:B:521:ARG:NH2	2.27	0.61
1:A:433:VAL:HG22	1:A:452:TYR:CD2	2.35	0.61
1:A:483:ASN:C	1:A:486:PRO:HD2	2.21	0.61
1:A:367:HIS:O	1:A:371:ALA:HB2	2.01	0.61
1:B:556:GLU:O	1:B:560:LYS:HG2	2.01	0.60
1:B:310:VAL:O	1:B:370:TYR:HE1	1.84	0.60
1:A:556:GLU:HG3	1:A:557:LYS:H	1.65	0.60
1:A:383:GLU:HB3	1:A:384:PRO:CD	2.31	0.60
1:B:372:LYS:O	1:B:375:ASP:HB2	2.00	0.60
1:B:57:GLU:OE1	1:B:57:GLU:HA	2.01	0.59
1:B:98:ARG:O	1:B:101:CYS:HB3	2.02	0.59
1:A:199:LYS:HG2	1:A:211:PHE:HE1	1.67	0.59
1:A:179:LEU:HB2	1:A:180:PRO:HD3	1.84	0.59
1:B:383:GLU:HB3	1:B:384:PRO:CD	2.26	0.59
1:B:485:ARG:HB3	1:B:486:PRO:HD3	1.84	0.59
1:B:511:ALA:HA	1:B:568:PHE:CE2	2.38	0.59
1:A:412:THR:O	1:A:416:PRO:HG3	2.02	0.59
1:A:305:LEU:HD21	1:A:333:GLU:HB3	1.84	0.59
1:A:240:LYS:O	1:A:244:GLU:HG3	2.03	0.59
1:A:87:MET:HE1	1:A:105:HIS:HB3	1.83	0.59
1:B:364:ALA:O	1:B:366:PRO:HD3	2.03	0.58
1:A:507:PHE:O	1:A:507:PHE:CD1	2.56	0.58
1:B:442:GLU:HA	1:B:445:ARG:HD2	1.85	0.58
1:B:120:VAL:HG21	1:B:175:ALA:CA	2.32	0.58
1:B:551:PHE:O	1:B:555:VAL:HG23	2.04	0.58
1:B:120:VAL:HG21	1:B:175:ALA:CB	2.34	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:214:TRP:CD1	1:B:343:VAL:HG11	2.39	0.57
1:A:417:GLN:CD	1:A:417:GLN:H	2.07	0.57
1:B:179:LEU:HB2	1:B:180:PRO:HD3	1.87	0.57
1:A:66:LEU:O	1:A:70:PHE:HD1	1.87	0.57
1:B:556:GLU:HG3	1:B:557:LYS:H	1.70	0.57
1:A:214:TRP:NE1	1:A:343:VAL:HG11	2.18	0.57
1:A:400:GLU:O	1:A:404:GLN:HG3	2.05	0.57
1:B:464:HIS:CE1	1:B:470:SER:H	2.23	0.56
1:B:212:LYS:O	1:B:216:VAL:HG23	2.05	0.56
1:B:464:HIS:HE1	1:B:470:SER:H	1.52	0.56
1:A:42:LEU:O	1:A:46:VAL:HG23	2.05	0.56
1:A:110:PRO:HB2	1:A:112:LEU:HG	1.88	0.56
1:B:531:GLU:HA	1:B:531:GLU:OE1	2.05	0.55
1:A:206:PHE:CE2	1:A:481:LEU:HD13	2.41	0.55
1:B:367:HIS:O	1:B:371:ALA:HB2	2.06	0.55
1:B:48:GLU:O	1:B:48:GLU:HG2	2.06	0.55
1:A:529:LEU:O	1:A:533:VAL:HG23	2.07	0.55
1:A:384:PRO:O	1:A:388:ILE:HG12	2.06	0.55
1:B:206:PHE:CD2	1:B:481:LEU:HD22	2.42	0.55
1:B:348:ARG:NH2	1:B:450:GLU:OE2	2.25	0.55
1:B:42:LEU:O	1:B:46:VAL:HG23	2.07	0.55
1:A:433:VAL:HG22	1:A:452:TYR:HD2	1.72	0.54
1:A:556:GLU:HG3	1:A:557:LYS:N	2.22	0.54
1:A:224:PRO:HD2	1:A:296:ASP:HB3	1.89	0.54
1:A:425:GLU:OE1	1:A:425:GLU:HA	2.07	0.54
1:A:578:ALA:O	1:A:581:ALA:HB3	2.07	0.54
1:A:106:LYS:HD3	1:A:147:PRO:HB2	1.90	0.54
1:B:344:VAL:O	1:B:348:ARG:HB2	2.09	0.53
1:A:19:PHE:O	1:A:23:VAL:HG23	2.08	0.53
1:B:542:GLU:O	1:B:545:LYS:HG2	2.10	0.53
1:B:373:VAL:HG13	1:B:374:PHE:HD1	1.74	0.52
1:B:543:GLN:O	1:B:546:ALA:HB3	2.09	0.52
1:B:50:ALA:O	1:B:54:VAL:HG23	2.09	0.52
1:B:75:CYS:HA	1:B:78:ALA:HB3	1.90	0.52
1:B:33:GLN:HE21	1:B:33:GLN:HA	1.73	0.52
1:B:483:ASN:C	1:B:486:PRO:HD2	2.30	0.52
1:A:49:PHE:O	1:A:49:PHE:HD2	1.93	0.52
1:A:543:GLN:O	1:A:547:VAL:HG23	2.10	0.52
1:B:404:GLN:NE2	1:B:428:ARG:HA	2.25	0.52
1:A:66:LEU:HB3	1:A:70:PHE:CE1	2.45	0.51
1:B:511:ALA:C	1:B:513:ILE:H	2.13	0.51
1:B:507:PHE:CD1	1:B:507:PHE:O	2.64	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:186:ARG:O	1:B:190:LYS:HG3	2.11	0.51
1:A:306:ALA:HA	1:A:310:VAL:HG22	1.93	0.51
1:B:153:GLU:O	1:B:157:PHE:HD1	1.93	0.51
1:B:389:LYS:O	1:B:393:GLU:HG3	2.11	0.51
1:B:196:GLN:HA	1:B:196:GLN:NE2	2.26	0.51
1:A:414:LYS:O	1:A:472:ARG:NH1	2.44	0.51
1:B:33:GLN:NE2	1:B:33:GLN:HA	2.26	0.50
1:A:509:PHE:O	1:A:568:PHE:CD1	2.64	0.50
1:A:571:GLU:N	1:A:571:GLU:OE1	2.43	0.50
1:A:38:ASP:OD1	1:A:38:ASP:N	2.44	0.50
1:B:505:GLU:O	1:B:506:THR:C	2.50	0.50
1:B:566:THR:O	1:B:570:GLU:N	2.40	0.50
1:B:405:ASN:O	1:B:409:VAL:HG23	2.11	0.50
1:B:344:VAL:HG21	1:B:450:GLU:HG2	1.93	0.50
1:B:492:GLU:OE1	1:B:492:GLU:HA	2.11	0.50
1:B:398:LEU:HB3	1:B:402:LYS:HB2	1.94	0.50
1:A:274:LYS:HD2	1:A:294:GLU:HG3	1.93	0.49
1:A:319:TYR:OH	1:A:358:GLU:HG2	2.12	0.49
1:A:310:VAL:HG23	1:A:311:GLU:N	2.27	0.49
1:B:36:PHE:O	1:B:40:VAL:HG23	2.12	0.49
1:A:420:THR:HB	1:A:421:PRO:HD3	1.94	0.49
1:B:417:GLN:H	1:B:417:GLN:CD	2.16	0.49
1:A:564:LYS:O	1:A:566:THR:N	2.45	0.49
1:A:87:MET:O	1:A:90:CYS:HB2	2.11	0.49
1:A:206:PHE:CE2	1:A:481:LEU:HB2	2.48	0.49
1:A:372:LYS:O	1:A:375:ASP:HB2	2.13	0.49
1:A:502:PHE:HE1	1:A:507:PHE:CE2	2.31	0.49
1:B:214:TRP:NE1	1:B:343:VAL:HG11	2.28	0.49
1:B:168:CYS:SG	1:B:177:CYS:C	2.91	0.48
1:B:556:GLU:HG3	1:B:557:LYS:N	2.28	0.48
1:B:120:VAL:HG21	1:B:175:ALA:HB2	1.95	0.48
1:A:566:THR:C	1:A:568:PHE:N	2.64	0.48
1:B:46:VAL:HG22	1:B:73:LYS:HE3	1.95	0.48
1:B:30:TYR:CE1	1:B:103:LEU:HD23	2.45	0.48
1:B:260:LEU:O	1:B:264:ILE:HG13	2.13	0.48
1:B:531:GLU:O	1:B:535:HIS:CD2	2.62	0.48
1:A:206:PHE:CD2	1:A:481:LEU:HB2	2.49	0.48
1:B:509:PHE:HB3	1:B:568:PHE:CD1	2.48	0.48
1:A:531:GLU:OE1	1:A:531:GLU:CA	2.59	0.48
1:B:458:ASN:O	1:B:462:VAL:HG22	2.13	0.48
1:B:425:GLU:OE1	1:B:425:GLU:HA	2.14	0.48
1:B:507:PHE:HZ	1:B:576:VAL:HG22	1.79	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:502:PHE:CG	1:A:503:ASN:N	2.81	0.47
1:B:25:ILE:O	1:B:29:GLN:HG3	2.14	0.47
1:B:283:LEU:HD12	1:B:283:LEU:C	2.34	0.47
1:A:115:LEU:HD22	1:A:145:ARG:NH1	2.29	0.47
1:B:400:GLU:O	1:B:404:GLN:HG3	2.13	0.47
1:B:26:ALA:HB2	1:B:250:LEU:HD12	1.96	0.47
1:A:153:GLU:O	1:A:157:PHE:HD1	1.97	0.47
1:A:373:VAL:HG13	1:A:374:PHE:HD1	1.79	0.47
1:B:9:HIS:O	1:B:13:ASP:HB2	2.14	0.47
1:B:19:PHE:O	1:B:23:VAL:HG23	2.14	0.47
1:B:206:PHE:HE2	1:B:481:LEU:HD13	1.77	0.47
1:B:383:GLU:OE2	1:B:387:LEU:CD1	2.61	0.47
1:A:117:ARG:HA	1:A:118:PRO:HD3	1.83	0.47
1:B:558:CYS:SG	1:B:568:PHE:N	2.88	0.47
1:B:178:LEU:O	1:B:179:LEU:C	2.53	0.47
1:B:282:PRO:HB2	1:B:285:GLU:OE1	2.14	0.47
1:A:66:LEU:HB3	1:A:70:PHE:CD1	2.50	0.46
1:A:417:GLN:OE1	1:A:417:GLN:N	2.28	0.46
1:B:33:GLN:CA	1:B:33:GLN:HE21	2.28	0.46
1:A:485:ARG:HB3	1:A:486:PRO:HD3	1.95	0.46
1:A:49:PHE:CD2	1:A:49:PHE:C	2.88	0.46
1:A:502:PHE:CE1	1:A:507:PHE:CE2	3.03	0.46
1:B:70:PHE:HD1	1:B:70:PHE:H	1.63	0.46
1:A:99:ASN:O	1:A:102:PHE:N	2.48	0.46
1:A:198:LEU:HA	1:A:458:ASN:ND2	2.31	0.46
1:B:290:ILE:O	1:B:293:VAL:HG12	2.15	0.46
1:A:566:THR:O	1:A:569:ALA:N	2.49	0.45
1:B:70:PHE:CD1	1:B:70:PHE:N	2.84	0.45
1:A:260:LEU:O	1:A:264:ILE:HG13	2.15	0.45
1:A:283:LEU:HG	1:A:284:LEU:HD23	1.98	0.45
1:B:38:ASP:O	1:B:42:LEU:HG	2.16	0.45
1:A:23:VAL:HG13	1:A:70:PHE:CE2	2.40	0.45
1:B:161:TYR:O	1:B:164:ALA:HB3	2.17	0.45
1:B:243:THR:HG22	1:B:247:HIS:HD2	1.82	0.45
1:B:420:THR:O	1:B:424:VAL:HG23	2.17	0.45
1:B:528:ALA:O	1:B:531:GLU:HB2	2.16	0.45
1:A:418:VAL:HB	1:A:423:LEU:HD21	2.00	0.45
1:B:141:GLU:O	1:B:144:ARG:HG2	2.17	0.45
1:B:119:GLU:OE1	1:B:119:GLU:HA	2.17	0.44
1:A:516:LEU:O	1:A:521:ARG:NH2	2.50	0.44
1:A:151:ALA:CB	1:A:152:PRO:HD3	2.42	0.44
1:A:370:TYR:CD1	1:A:370:TYR:C	2.91	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:49:PHE:CE2	1:A:53:CYS:SG	3.11	0.44
1:A:429:ASN:O	1:A:432:LYS:N	2.49	0.44
1:B:14:LEU:CD1	1:B:251:LEU:HD22	2.47	0.44
1:B:530:VAL:HG12	1:B:534:LYS:HD2	2.00	0.44
1:A:344:VAL:HG21	1:A:450:GLU:HG2	2.00	0.44
1:A:528:ALA:O	1:A:531:GLU:HB2	2.17	0.44
1:B:70:PHE:HD1	1:B:70:PHE:N	2.15	0.44
1:B:525:LYS:O	1:B:528:ALA:HB3	2.17	0.44
1:B:139:LEU:HD21	1:B:158:ALA:HB2	1.99	0.44
1:B:173:ASP:OD2	1:B:176:ALA:HB2	2.18	0.44
1:A:509:PHE:O	1:A:510:HIS:C	2.56	0.44
1:B:283:LEU:HG	1:B:284:LEU:HD23	2.00	0.44
1:A:120:VAL:HG11	1:A:174:LYS:CB	2.48	0.43
1:A:511:ALA:C	1:A:513:ILE:H	2.22	0.43
1:B:563:ASP:OD1	1:B:567:CYS:SG	2.76	0.43
1:B:408:LEU:HD12	1:B:526:GLN:HG2	2.01	0.43
1:A:34:CYS:HA	1:A:35:PRO:HD3	1.86	0.43
1:B:359:LYS:CG	1:B:360:CYS:N	2.80	0.43
1:B:221:GLN:O	1:B:295:ASN:HB3	2.18	0.43
1:B:384:PRO:O	1:B:388:ILE:HG12	2.18	0.43
1:B:7:VAL:HG13	1:B:8:ALA:N	2.34	0.43
1:A:30:TYR:HE1	1:A:103:LEU:HD23	1.83	0.43
1:B:215:ALA:HB3	1:B:235:VAL:HG13	2.00	0.43
1:A:168:CYS:SG	1:A:177:CYS:O	2.76	0.43
1:A:333:GLU:OE1	1:A:337:ARG:NH2	2.50	0.43
1:A:290:ILE:O	1:A:293:VAL:HG12	2.19	0.43
1:A:549:ASP:O	1:A:552:ALA:HB3	2.18	0.43
1:A:272:SER:HB3	1:A:275:LEU:HG	2.00	0.43
1:B:367:HIS:HA	1:B:370:TYR:CZ	2.53	0.43
1:A:564:LYS:O	1:A:565:GLU:C	2.57	0.43
1:A:231:VAL:O	1:A:234:LEU:N	2.51	0.43
1:A:420:THR:HG23	1:A:530:VAL:HG11	2.01	0.43
1:A:453:LEU:HD23	1:A:485:ARG:HG3	2.01	0.43
1:A:67:HIS:CE1	1:A:99:ASN:ND2	2.87	0.43
1:B:349:LEU:HD22	1:B:377:PHE:CG	2.54	0.43
1:A:23:VAL:O	1:A:27:PHE:HD1	2.02	0.42
1:A:120:VAL:HG21	1:A:175:ALA:CA	2.45	0.42
1:A:178:LEU:O	1:A:179:LEU:C	2.57	0.42
1:A:441:PRO:O	1:A:442:GLU:C	2.57	0.42
1:A:153:GLU:O	1:A:156:PHE:HB3	2.19	0.42
1:A:331:LEU:HD21	1:A:347:LEU:HD23	2.00	0.42
1:A:151:ALA:HB3	1:A:152:PRO:CD	2.42	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:558:CYS:SG	1:A:568:PHE:CD2	3.12	0.42
1:A:498:VAL:HG13	1:A:498:VAL:O	2.19	0.42
1:A:14:LEU:HD13	1:A:22:LEU:HD12	2.00	0.42
1:B:441:PRO:O	1:B:443:ALA:N	2.52	0.42
1:B:404:GLN:HE21	1:B:428:ARG:HA	1.83	0.42
1:A:113:PRO:O	1:A:114:ARG:C	2.56	0.42
1:A:109:ASN:HD22	1:A:109:ASN:HA	1.54	0.42
1:B:95:GLU:OE1	1:B:99:ASN:HB2	2.19	0.42
1:A:224:PRO:CB	1:A:299:PRO:HD3	2.43	0.42
1:A:464:HIS:CE1	1:A:469:VAL:H	2.38	0.42
1:A:567:CYS:O	1:A:571:GLU:N	2.52	0.42
1:A:344:VAL:O	1:A:348:ARG:HB2	2.20	0.42
1:A:99:ASN:O	1:A:100:GLU:C	2.58	0.42
1:A:429:ASN:O	1:A:431:GLY:N	2.53	0.42
1:A:117:ARG:HB2	1:A:123:MET:CE	2.50	0.42
1:B:187:ASP:O	1:B:188:GLU:C	2.58	0.42
1:A:310:VAL:O	1:A:370:TYR:HE1	2.03	0.42
1:B:420:THR:HG23	1:B:530:VAL:HG11	2.02	0.41
1:B:275:LEU:O	1:B:276:LYS:C	2.58	0.41
1:A:511:ALA:O	1:A:513:ILE:N	2.53	0.41
1:A:558:CYS:SG	1:A:568:PHE:N	2.93	0.41
1:A:555:VAL:O	1:A:559:CYS:HB2	2.19	0.41
1:A:39:HIS:O	1:A:43:VAL:HG23	2.20	0.41
1:A:437:CYS:SG	1:A:448:CYS:C	2.98	0.41
1:B:532:LEU:HD11	1:B:547:VAL:HG11	2.02	0.41
1:B:518:GLU:O	1:B:521:ARG:HB3	2.20	0.41
1:B:370:TYR:C	1:B:370:TYR:CD1	2.93	0.41
1:B:532:LEU:HA	1:B:532:LEU:HD23	1.96	0.41
1:A:403:PHE:CE2	1:A:407:LEU:HD11	2.55	0.41
1:B:521:ARG:HB2	1:B:521:ARG:HH11	1.85	0.41
1:B:311:GLU:O	1:B:312:SER:C	2.58	0.41
1:A:26:ALA:HB2	1:A:250:LEU:HD12	2.01	0.41
1:B:179:LEU:CB	1:B:180:PRO:HD3	2.51	0.41
1:A:25:ILE:HD13	1:A:154:LEU:HD23	2.03	0.41
1:B:168:CYS:SG	1:B:178:LEU:N	2.93	0.41
1:B:187:ASP:O	1:B:190:LYS:N	2.53	0.41
1:B:310:VAL:HG21	1:B:374:PHE:CE1	2.56	0.41
1:A:179:LEU:H	1:A:179:LEU:HD23	1.84	0.41
1:B:48:GLU:O	1:B:48:GLU:CG	2.69	0.41
1:A:9:HIS:ND1	1:A:13:ASP:OD2	2.53	0.41
1:B:57:GLU:O	1:B:59:ALA:N	2.54	0.40
1:A:49:PHE:HD2	1:A:49:PHE:C	2.25	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:232:SER:O	1:A:236:THR:OG1	2.36	0.40
1:B:453:LEU:HD13	2:B:2001:DZP:H203	2.02	0.40
1:A:508:THR:CG2	1:A:509:PHE:N	2.84	0.40
1:B:141:GLU:OE1	1:B:144:ARG:HD3	2.21	0.40
1:B:14:LEU:HD13	1:B:22:LEU:HD12	2.04	0.40
1:A:507:PHE:C	1:A:507:PHE:CD1	2.95	0.40
1:A:511:ALA:HA	1:A:568:PHE:CE2	2.56	0.40
1:B:57:GLU:HB3	1:B:58:SER:H	1.47	0.40
1:A:141:GLU:OE1	1:A:141:GLU:HA	2.22	0.40
1:A:330:PHE:HE2	1:A:350:ALA:HA	1.86	0.40
1:A:218:ARG:HH21	1:A:222:ARG:HH21	1.70	0.40
1:B:21:ALA:O	1:B:25:ILE:HG13	2.22	0.40
1:B:494:ASP:OD1	1:B:494:ASP:C	2.59	0.40
1:A:351:LYS:HA	1:A:351:LYS:HD2	1.87	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	576/585 (98%)	481 (84%)	78 (14%)	17 (3%)	7	32
1	B	576/585 (98%)	490 (85%)	66 (12%)	20 (4%)	6	27
All	All	1152/1170 (98%)	971 (84%)	144 (12%)	37 (3%)	6	30

All (37) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	59	ALA
1	A	300	ALA
1	A	565	GLU
1	B	54	VAL
1	B	57	GLU
1	B	58	SER

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Mol	Chain	Res	Type
1	B	60	GLU
1	B	129	ASP
1	A	17	GLU
1	A	58	SER
1	B	85	GLY
1	B	114	ARG
1	B	276	LYS
1	B	565	GLU
1	A	312	SER
1	A	367	HIS
1	A	512	ASP
1	B	363	ALA
1	B	367	HIS
1	B	442	GLU
1	A	430	LEU
1	B	64	LYS
1	B	557	LYS
1	A	114	ARG
1	A	179	LEU
1	A	180	PRO
1	A	255	ASP
1	B	77	VAL
1	B	379	PRO
1	B	380	LEU
1	A	366	PRO
1	A	390	GLN
1	B	118	PRO
1	A	241	VAL
1	A	315	VAL
1	B	179	LEU
1	B	315	VAL

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	421/511 (82%)	403 (96%)	18 (4%)	40 81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	424/511 (83%)	406 (96%)	18 (4%)	40	82
All	All	845/1022 (83%)	809 (96%)	36 (4%)	40	81

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	HIS
1	A	37	GLU
1	A	38	ASP
1	A	49	PHE
1	A	66	LEU
1	A	72	ASP
1	A	79	THR
1	A	109	ASN
1	A	179	LEU
1	A	186	ARG
1	A	222	ARG
1	A	249	ASP
1	A	268	GLN
1	A	334	TYR
1	A	435	SER
1	A	467	THR
1	A	543	GLN
1	A	550	ASP
1	B	34	CYS
1	B	38	ASP
1	B	57	GLU
1	B	79	THR
1	B	105	HIS
1	B	179	LEU
1	B	245	CYS
1	B	283	LEU
1	B	292	GLU
1	B	334	TYR
1	B	352	THR
1	B	375	ASP
1	B	408	LEU
1	B	415	VAL
1	B	419	SER
1	B	459	GLN
1	B	566	THR
1	B	568	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	ASN
1	A	196	GLN
1	A	338	HIS
1	A	367	HIS
1	A	464	HIS
1	A	483	ASN
1	B	33	GLN
1	B	99	ASN
1	B	196	GLN
1	B	247	HIS
1	B	385	GLN
1	B	404	GLN
1	B	464	HIS
1	B	483	ASN
1	B	535	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	DZP	A	2001	-	22,22,22	1.55	5 (22%)	31,31,31	0.72	0
2	DZP	B	2001	-	22,22,22	1.51	4 (18%)	31,31,31	0.75	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	DZP	A	2001	-	-	0/4/21/21	0/1/3/3
2	DZP	B	2001	-	-	0/4/21/21	0/1/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2001	DZP	C4-C12	2.88	1.44	1.39
2	A	2001	DZP	C4-C12	2.61	1.43	1.39
2	B	2001	DZP	C5-N15	2.54	1.32	1.28
2	A	2001	DZP	C11-C12	2.44	1.43	1.39
2	A	2001	DZP	C13-C6	2.43	1.45	1.41
2	B	2001	DZP	C18-N2	2.37	1.41	1.37
2	B	2001	DZP	C13-C6	2.14	1.44	1.41
2	A	2001	DZP	C5-N15	2.11	1.31	1.28
2	A	2001	DZP	C12-C5	2.10	1.52	1.49

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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, 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	578/585 (98%)	-0.17	22 (3%)	38 18	33, 77, 141, 154	0
1	B	578/585 (98%)	-0.15	29 (5%)	28 14	35, 80, 145, 159	0
All	All	1156/1170 (98%)	-0.16	51 (4%)	33 16	33, 78, 144, 159	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	79	THR	5.1
1	B	55	ALA	4.9
1	A	572	GLY	4.8
1	B	90	CYS	4.5
1	B	576	VAL	4.3
1	B	91	CYS	3.8
1	A	562	ASP	3.7
1	B	564	LYS	3.7
1	A	561	ALA	3.5
1	B	562	ASP	3.5
1	B	300	ALA	3.4
1	B	569	ALA	3.3
1	B	83	THR	3.3
1	B	515	THR	3.3
1	A	558	CYS	3.2
1	A	56	ASP	3.1
1	B	79	THR	3.0
1	B	572	GLY	3.0
1	A	83	THR	3.0
1	A	78	ALA	2.9
1	B	577	ALA	2.9
1	A	564	LYS	2.9
1	A	301	ASP	2.9
1	B	364	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	512	ASP	2.7
1	A	300	ALA	2.7
1	A	563	ASP	2.7
1	B	573	LYS	2.6
1	A	577	ALA	2.6
1	B	82	GLU	2.6
1	A	573	LYS	2.5
1	B	566	THR	2.4
1	B	56	ASP	2.4
1	B	560	LYS	2.3
1	A	91	CYS	2.3
1	A	506	THR	2.3
1	B	558	CYS	2.3
1	A	92	ALA	2.2
1	B	50	ALA	2.2
1	B	78	ALA	2.2
1	B	467	THR	2.1
1	B	561	ALA	2.1
1	B	510	HIS	2.1
1	A	576	VAL	2.1
1	A	560	LYS	2.1
1	B	508	THR	2.1
1	A	510	HIS	2.1
1	B	80	LEU	2.1
1	A	26	ALA	2.1
1	B	511	ALA	2.0
1	A	27	PHE	2.0

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 carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	DZP	B	2001	20/20	0.22	1.56	57,58,63,63	0
2	DZP	A	2001	20/20	0.22	1.23	50,56,60,62	0

6.5 Other polymers ⓘ

There are no such residues in this entry.