Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 06:51 PM EDT

PDB ID	:	1EW2
Title	:	CRYSTAL STRUCTURE OF A HUMAN PHOSPHATASE
Authors	:	Le Du, M.H.; Stigbrand, T.; Taussig, M.J.; Menez, A.; Stura, E.A.
Deposited on	:	2000-04-21
Resolution	:	1.82 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 (i) 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 (i)) were used in the production of this report:

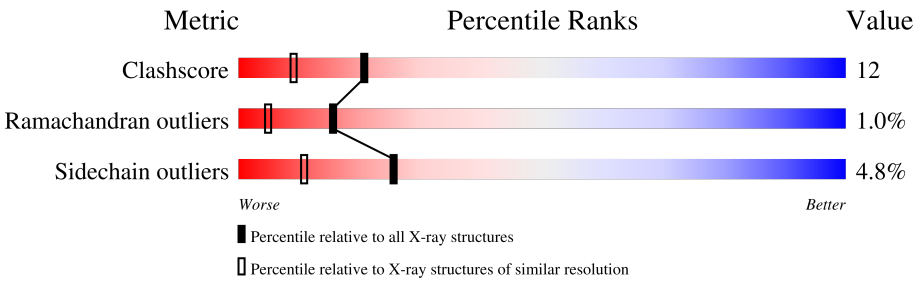
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 1.82 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))
Clashscore	141614	8401 (1.84-1.80)
Ramachandran outliers	138981	8290 (1.84-1.80)
Sidechain outliers	138945	8290 (1.84-1.80)

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	513	71% 17% • • 7%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	PO4	A	1005	-	X	X	-

2 Entry composition (i)

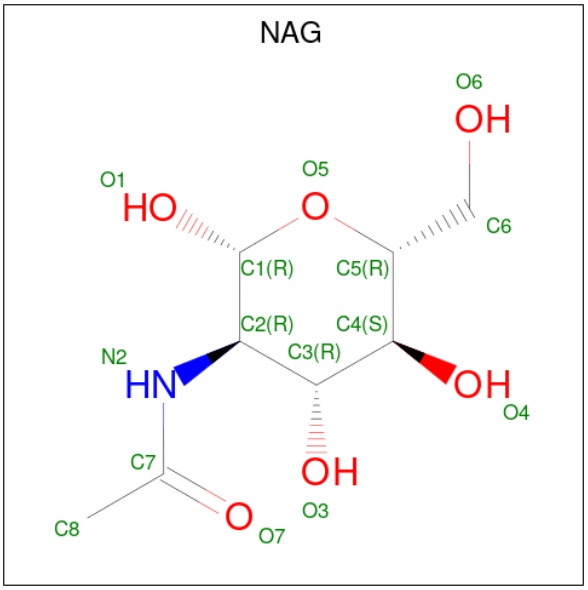
There are 6 unique types of molecules in this entry. The entry contains 4277 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PHOSPHATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	479	3652	2285	651	698	18	0	0	0

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	14	8	1	5	0	0

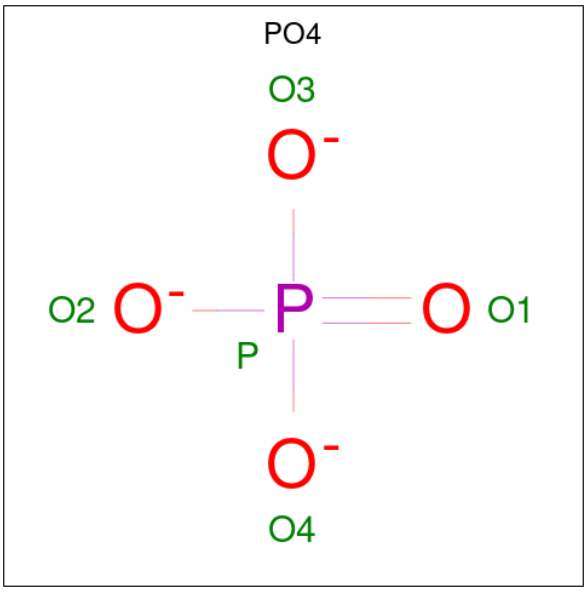
- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
3	A	2	2	2	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total Mg 2 2	0	0

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total O P 5 4 1	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	602	Total O 602 602	0	0

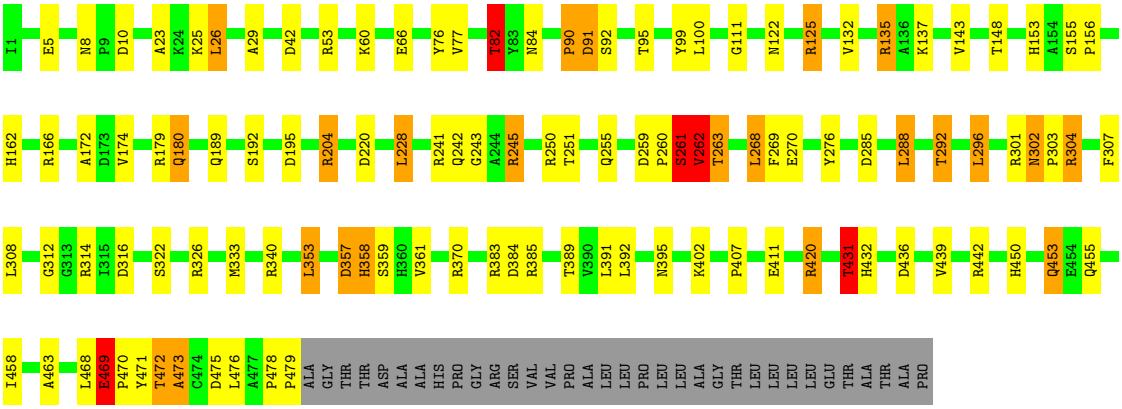
3 Residue-property plots (i)

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. 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. 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.

Note EDS was not executed.

• Molecule 1: PHOSPHATASE

Chain A:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	88.80Å 114.50Å 106.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	22.00 – 1.82	Depositor
% Data completeness (in resolution range)	(Not available) (22.00-1.82)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.185 , 0.242	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4277	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, PO4, NAG, ZN

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	0.71	1/3733 (0.0%)	1.63	67/5066 (1.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	479	PRO	N-CD	5.52	1.55	1.47

All (67) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	357	ASP	CB-CG-OD1	18.84	135.26	118.30
1	A	472	THR	N-CA-CB	18.72	145.87	110.30
1	A	357	ASP	CB-CG-OD2	-15.14	104.67	118.30
1	A	42	ASP	CB-CG-OD2	15.04	131.84	118.30
1	A	125	ARG	CD-NE-CZ	14.51	143.91	123.60
1	A	125	ARG	NE-CZ-NH1	12.13	126.36	120.30
1	A	135	ARG	NE-CZ-NH1	-11.74	114.43	120.30
1	A	301	ARG	CD-NE-CZ	10.59	138.43	123.60
1	A	262	VAL	CA-C-N	10.19	139.62	117.20
1	A	472	THR	CA-CB-CG2	9.39	125.55	112.40
1	A	301	ARG	NE-CZ-NH1	8.99	124.80	120.30
1	A	262	VAL	O-C-N	-8.90	108.46	122.70

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	245	ARG	NE-CZ-NH2	-8.51	116.04	120.30
1	A	431	THR	N-CA-CB	-8.48	94.19	110.30
1	A	420	ARG	NE-CZ-NH1	8.30	124.45	120.30
1	A	82	THR	CB-CA-C	-8.16	89.57	111.60
1	A	82	THR	CA-CB-OG1	7.82	125.41	109.00
1	A	314	ARG	NE-CZ-NH1	-7.67	116.47	120.30
1	A	316	ASP	CB-CG-OD1	-7.59	111.47	118.30
1	A	316	ASP	CB-CG-OD2	7.46	125.02	118.30
1	A	204	ARG	NE-CZ-NH2	-7.39	116.61	120.30
1	A	60	LYS	CA-C-N	7.37	133.41	117.20
1	A	166	ARG	NE-CZ-NH1	-7.20	116.70	120.30
1	A	370	ARG	NE-CZ-NH2	7.16	123.88	120.30
1	A	53	ARG	NE-CZ-NH2	-6.93	116.83	120.30
1	A	370	ARG	NE-CZ-NH1	-6.88	116.86	120.30
1	A	261	SER	N-CA-CB	6.80	120.70	110.50
1	A	245	ARG	NE-CZ-NH1	6.78	123.69	120.30
1	A	180	GLN	CB-CG-CD	6.74	129.12	111.60
1	A	385	ARG	NE-CZ-NH1	6.66	123.63	120.30
1	A	384	ASP	CB-CG-OD2	-6.56	112.39	118.30
1	A	76	TYR	CB-CG-CD1	-6.53	117.08	121.00
1	A	285	ASP	CB-CG-OD1	6.34	124.01	118.30
1	A	276	TYR	CB-CG-CD2	-6.33	117.20	121.00
1	A	180	GLN	CA-CB-CG	6.13	126.88	113.40
1	A	353	LEU	CA-CB-CG	6.09	129.32	115.30
1	A	60	LYS	O-C-N	-5.98	113.14	122.70
1	A	263	THR	N-CA-CB	5.97	121.64	110.30
1	A	384	ASP	CB-CG-OD1	5.95	123.66	118.30
1	A	472	THR	CA-CB-OG1	-5.95	96.50	109.00
1	A	472	THR	N-CA-C	-5.85	95.20	111.00
1	A	326	ARG	NE-CZ-NH2	5.83	123.21	120.30
1	A	192	SER	N-CA-CB	-5.80	101.80	110.50
1	A	135	ARG	NE-CZ-NH2	5.79	123.19	120.30
1	A	250	ARG	NE-CZ-NH1	-5.72	117.44	120.30
1	A	66	GLU	OE1-CD-OE2	-5.71	116.45	123.30
1	A	42	ASP	OD1-CG-OD2	-5.64	112.58	123.30
1	A	250	ARG	NE-CZ-NH2	5.61	123.11	120.30
1	A	220	ASP	CB-CG-OD2	5.58	123.32	118.30
1	A	358	HIS	CA-CB-CG	-5.52	104.21	113.60
1	A	453	GLN	CB-CG-CD	5.49	125.86	111.60
1	A	436	ASP	CB-CG-OD2	-5.44	113.40	118.30
1	A	471	TYR	C-N-CA	-5.44	108.11	121.70
1	A	270	GLU	OE1-CD-OE2	-5.41	116.81	123.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	241	ARG	NE-CZ-NH2	5.34	122.97	120.30
1	A	312	GLY	N-CA-C	-5.26	99.94	113.10
1	A	316	ASP	N-CA-CB	-5.25	101.16	110.60
1	A	135	ARG	CA-CB-CG	-5.21	101.93	113.40
1	A	285	ASP	CB-CG-OD2	-5.18	113.64	118.30
1	A	340	ARG	NE-CZ-NH1	5.18	122.89	120.30
1	A	304	ARG	NE-CZ-NH2	-5.17	117.72	120.30
1	A	301	ARG	NH1-CZ-NH2	-5.16	113.73	119.40
1	A	358	HIS	ND1-CE1-NE2	5.15	121.23	109.90
1	A	269	PHE	CA-C-N	5.13	128.50	117.20
1	A	478	PRO	C-N-CD	5.11	139.13	128.40
1	A	469	GLU	CB-CA-C	5.10	120.61	110.40
1	A	442	ARG	N-CA-CB	5.09	119.76	110.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	468	LEU	Mainchain
1	A	82	THR	Mainchain

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	3652	0	3539	89	0
2	A	14	0	13	1	0
3	A	2	0	0	0	0
4	A	2	0	0	0	0
5	A	5	0	0	2	0
6	A	602	0	0	28	4
All	All	4277	0	3552	89	4

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 12.

All (89) 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:26:LEU:HD21	6:A:1556:HOH:O	1.16	1.32
1:A:23:ALA:O	1:A:26:LEU:HD22	1.71	0.89
1:A:23:ALA:O	1:A:26:LEU:CD2	2.26	0.84
1:A:162:HIS:H	1:A:189:GLN:HE22	1.23	0.83
1:A:135:ARG:HD2	1:A:469:GLU:HA	1.64	0.78
1:A:383:ARG:HG2	6:A:1328:HOH:O	1.87	0.73
1:A:82:THR:HG23	6:A:1576:HOH:O	1.88	0.73
1:A:402:LYS:HE3	1:A:407:PRO:HB3	1.70	0.72
1:A:29:ALA:HB2	6:A:1603:HOH:O	1.90	0.72
1:A:135:ARG:HH11	1:A:135:ARG:HG3	1.59	0.66
1:A:259:ASP:OD2	1:A:260:PRO:O	2.14	0.66
1:A:135:ARG:HD3	6:A:1605:HOH:O	1.96	0.65
1:A:469:GLU:HA	6:A:1605:HOH:O	1.96	0.64
1:A:82:THR:CB	6:A:1576:HOH:O	2.44	0.63
1:A:82:THR:CG2	6:A:1576:HOH:O	2.46	0.62
1:A:472:THR:O	1:A:473:ALA:HB3	2.01	0.61
1:A:82:THR:CG2	1:A:90:PRO:HG3	2.31	0.60
1:A:472:THR:OG1	1:A:473:ALA:N	2.28	0.60
1:A:453:GLN:HG2	6:A:1046:HOH:O	2.02	0.60
1:A:469:GLU:CB	1:A:470:PRO:CD	2.80	0.59
1:A:26:LEU:CD2	6:A:1556:HOH:O	2.00	0.58
1:A:135:ARG:HG3	1:A:135:ARG:NH1	2.19	0.58
1:A:10:ASP:HB2	6:A:1558:HOH:O	2.04	0.58
1:A:92:SER:OG	5:A:1005:PO4:P	2.62	0.58
1:A:137:LYS:HE2	1:A:195:ASP:HB3	1.86	0.57
1:A:431:THR:HG23	1:A:432:HIS:O	2.04	0.57
1:A:95:THR:HG22	1:A:455:GLN:HE22	1.71	0.56
1:A:5:GLU:HA	6:A:1469:HOH:O	2.05	0.55
1:A:472:THR:HG22	6:A:1379:HOH:O	2.05	0.55
1:A:302:ASN:ND2	1:A:304:ARG:H	2.04	0.55
1:A:353:LEU:HD12	1:A:458:ILE:HG23	1.88	0.54
1:A:204:ARG:HG2	1:A:268:LEU:HB3	1.88	0.54
1:A:135:ARG:NH1	6:A:1132:HOH:O	2.24	0.54
1:A:172:ALA:HA	1:A:228:LEU:HD22	1.90	0.54
1:A:135:ARG:CD	1:A:469:GLU:HA	2.38	0.53
1:A:242:GLN:OE1	6:A:1601:HOH:O	0.53	0.53
1:A:82:THR:HG21	1:A:90:PRO:HG3	1.90	0.53
1:A:99:TYR:CD1	1:A:100:LEU:HG	2.44	0.53
1:A:243:GLY:O	1:A:262:VAL:O	2.28	0.52
1:A:153:HIS:HD2	6:A:1115:HOH:O	1.90	0.52
1:A:135:ARG:HH11	1:A:135:ARG:CG	2.18	0.52
1:A:162:HIS:H	1:A:189:GLN:NE2	2.02	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:260:PRO:O	1:A:261:SER:CB	2.57	0.52
1:A:389:THR:O	1:A:392:LEU:HD21	2.10	0.52
1:A:472:THR:O	1:A:473:ALA:CB	2.58	0.52
1:A:469:GLU:CA	6:A:1605:HOH:O	2.55	0.52
1:A:439:VAL:HG21	1:A:458:ILE:HD11	1.92	0.51
1:A:26:LEU:HD23	6:A:1346:HOH:O	2.09	0.51
1:A:162:HIS:HE1	6:A:1079:HOH:O	1.93	0.51
1:A:82:THR:HG21	1:A:358:HIS:HB3	1.92	0.50
1:A:251:THR:O	1:A:255:GLN:HG3	2.10	0.50
1:A:111:GLY:O	1:A:162:HIS:HD2	1.94	0.50
1:A:411:GLU:HG2	6:A:1566:HOH:O	2.12	0.49
1:A:288:LEU:O	1:A:292:THR:HG23	2.12	0.48
1:A:251:THR:HG23	6:A:1315:HOH:O	2.14	0.48
1:A:463:ALA:HB1	1:A:469:GLU:CB	2.44	0.48
1:A:358:HIS:HD2	1:A:359:SER:O	1.97	0.47
1:A:383:ARG:HD2	1:A:383:ARG:O	2.15	0.47
1:A:391:LEU:C	1:A:392:LEU:HD22	2.34	0.47
1:A:395:ASN:HD21	1:A:431:THR:CG2	2.27	0.47
1:A:135:ARG:CD	6:A:1605:HOH:O	2.58	0.46
1:A:125:ARG:HD3	6:A:1523:HOH:O	2.16	0.46
1:A:450:HIS:HE1	6:A:1228:HOH:O	1.99	0.46
1:A:439:VAL:CG2	1:A:458:ILE:HD11	2.46	0.45
1:A:92:SER:HG	5:A:1005:PO4:P	2.40	0.45
1:A:82:THR:CG2	1:A:357:ASP:O	2.65	0.44
1:A:143:VAL:HG12	1:A:307:PHE:HB3	1.98	0.44
1:A:302:ASN:HD22	1:A:303:PRO:N	2.15	0.43
1:A:304:ARG:HG2	6:A:1506:HOH:O	2.17	0.43
1:A:302:ASN:HD22	1:A:304:ARG:H	1.63	0.43
1:A:333:MET:HE2	6:A:1459:HOH:O	2.18	0.43
1:A:77:VAL:HA	1:A:439:VAL:O	2.18	0.42
1:A:135:ARG:HD2	1:A:469:GLU:CA	2.43	0.42
1:A:262:VAL:O	1:A:263:THR:CB	2.67	0.42
1:A:395:ASN:HD21	1:A:431:THR:HG21	1.84	0.42
1:A:389:THR:O	1:A:392:LEU:CD2	2.67	0.42
1:A:84:ASN:HD21	1:A:91:ASP:H	1.66	0.42
1:A:259:ASP:HA	1:A:260:PRO:HD2	1.94	0.42
1:A:260:PRO:O	1:A:261:SER:HB3	2.20	0.41
1:A:296:LEU:HD13	1:A:308:LEU:HD22	2.03	0.41
1:A:302:ASN:HD22	1:A:302:ASN:C	2.23	0.41
1:A:469:GLU:CB	1:A:470:PRO:HD2	2.49	0.41
1:A:155:SER:N	1:A:156:PRO:HD2	2.35	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:HIS:CD2	6:A:1115:HOH:O	2.71	0.41
1:A:148:THR:HG23	1:A:288:LEU:HG	2.01	0.41
1:A:392:LEU:HD23	6:A:1051:HOH:O	2.20	0.41
1:A:322:SER:O	1:A:420:ARG:HA	2.21	0.41
1:A:174:VAL:O	1:A:179:ARG:NH1	2.53	0.40
1:A:122:ASN:O	2:A:1000:NAG:H82	2.20	0.40

All (4) 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)
6:A:1598:HOH:O	6:A:1598:HOH:O[3_655]	0.80	1.40
6:A:1133:HOH:O	6:A:1222:HOH:O[4_555]	1.91	0.29
6:A:1035:HOH:O	6:A:1399:HOH:O[3_655]	2.12	0.08
6:A:1133:HOH:O	6:A:1137:HOH:O[4_555]	2.19	0.01

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	477/513 (93%)	460 (96%)	12 (2%)	5 (1%)	15 5

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	469	GLU
1	A	261	SER
1	A	473	ALA
1	A	262	VAL
1	A	361	VAL

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	377/408 (92%)	359 (95%)	18 (5%)	25 11

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	25	LYS
1	A	26	LEU
1	A	82	THR
1	A	90	PRO
1	A	91	ASP
1	A	132	VAL
1	A	180	GLN
1	A	228	LEU
1	A	245	ARG
1	A	268	LEU
1	A	288	LEU
1	A	292	THR
1	A	296	LEU
1	A	302	ASN
1	A	431	THR
1	A	475	ASP
1	A	476	LEU

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

Mol	Chain	Res	Type
1	A	8	ASN
1	A	84	ASN
1	A	127	ASN
1	A	134	ASN
1	A	153	HIS
1	A	162	HIS
1	A	184	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	189	GLN
1	A	193	ASN
1	A	302	ASN
1	A	358	HIS
1	A	445	GLN
1	A	450	HIS
1	A	455	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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	NAG	A	1000	1	14,14,15	1.62	4 (28%)	17,19,21	2.23	3 (17%)
5	PO4	A	1005	3	4,4,4	2.79	3 (75%)	6,6,6	3.60	5 (83%)

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	NAG	A	1000	1	-	2/6/23/26	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1005	PO4	P-O4	-3.60	1.43	1.54
2	A	1000	NAG	O5-C1	3.24	1.48	1.43
5	A	1005	PO4	P-O2	-3.03	1.45	1.54
2	A	1000	NAG	O7-C7	-2.99	1.16	1.23
5	A	1005	PO4	P-O1	2.95	1.57	1.50
2	A	1000	NAG	C1-C2	-2.40	1.48	1.52
2	A	1000	NAG	C2-N2	2.25	1.50	1.46

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1000	NAG	C2-N2-C7	7.50	133.59	122.90
5	A	1005	PO4	O4-P-O3	6.06	127.42	107.97
5	A	1005	PO4	O4-P-O2	3.82	120.22	107.97
5	A	1005	PO4	O4-P-O1	-3.56	97.86	110.89
2	A	1000	NAG	C4-C3-C2	-3.04	106.56	111.02
5	A	1005	PO4	O3-P-O2	-2.68	99.36	107.97
5	A	1005	PO4	O3-P-O1	-2.28	102.56	110.89
2	A	1000	NAG	O7-C7-C8	2.20	126.13	122.06

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1000	NAG	C8-C7-N2-C2
2	A	1000	NAG	O7-C7-N2-C2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1000	NAG	1	0
5	A	1005	PO4	2	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 (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.