



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:55 PM GMT

PDB ID : 1D6V
Title : CONFORMATION EFFECTS IN BIOLOGICAL CATALYSIS INTRODUCED BY OXY-COPE ANTIBODY MATURATION
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Deposited on : 1999-10-15
Resolution : 2.00 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

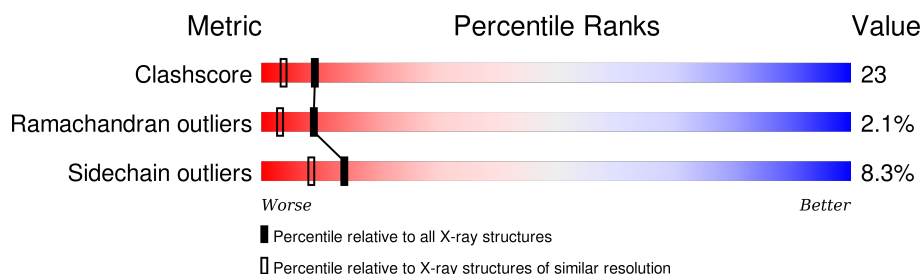
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	211	 68% 27% . .
2	H	221	 59% 33% 8% .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3486 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 CHIMERIC GERMLINE PRECURSOR OF OXY-COPE CATALYTIC ANTIBODY AZ-28 (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	211	Total	C	N	O	S	0	0	0
			1652	1038	272	335	7			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	34	SER	ASN	SEE REMARK 999	UNP P01834
L	51	ALA	THR	SEE REMARK 999	UNP P01834
L	108	ARG	-	NOT IN SWS ENTRY	UNP P01834

- Molecule 2 is a protein called CHIMERIC GERMLINE PRECURSOR OF OXY-COPE CATALYTIC ANTIBODY AZ-28 (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	221	Total	C	N	O	S	0	0	0
			1664	1053	271	334	6			

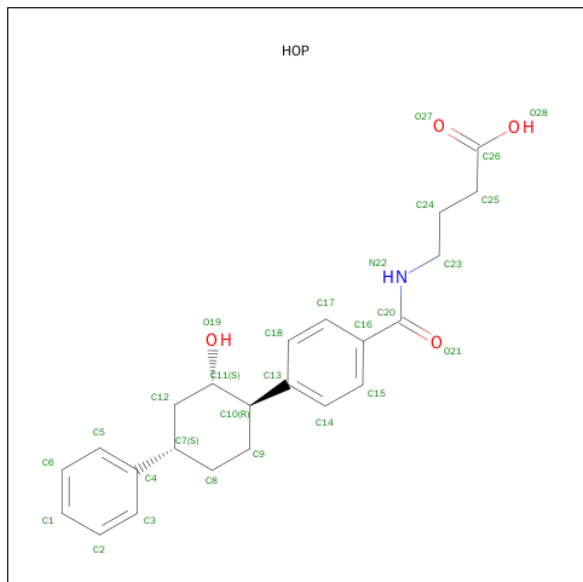
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	32	TYR	PHE	SEE REMARK 999	UNP P01857
H	56	SER	GLY	SEE REMARK 999	UNP P01857
H	58	ASN	HIS	SEE REMARK 999	UNP P01857
H	73	THR	LYS	SEE REMARK 999	UNP P01857

- Molecule 3 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	4	Total	Cd	0	0
			4	4		

- Molecule 4 is (1S,2S,5S)2-(4-GLUTARIDYLBENZYL)-5-PHENYL-1-CYCLOHEXANOL (three-letter code: HOP) (formula: C₂₃H₂₇NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	H	1	Total	C	N	O	0	0
			28	23	1	4		

- Molecule 5 is water.

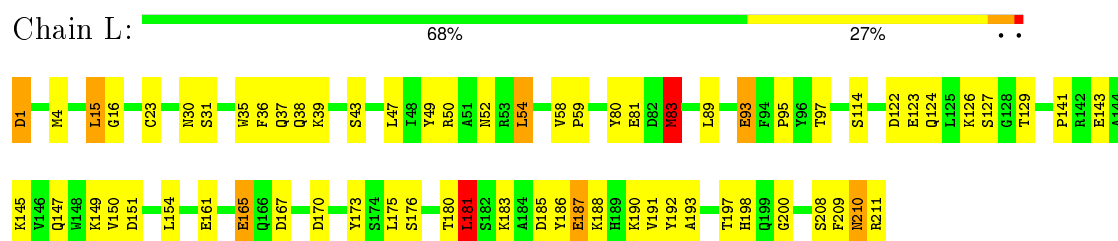
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	H	56	Total	O	0	0
			56	56		
5	L	82	Total	O	0	0
			82	82		

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.

Note EDS was not executed.

- Molecule 1: CHIMERIC GERMLINE PRECURSOR OF OXY-COPE CATALYTIC ANTI-BODY AZ-28 (LIGHT CHAIN)



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.72Å 64.45Å 73.87Å 90.00° 99.09° 90.00°	Depositor
Resolution (Å)	9.92 – 2.00	Depositor
% Data completeness (in resolution range)	83.1 (9.92-2.00)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.4	Depositor
R, R_{free}	0.224 , 0.264	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3486	wwPDB-VP
Average B, all atoms (Å ²)	36.0	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: HOP, CD

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	L	1.04	1/1688 (0.1%)	1.06	3/2284 (0.1%)
2	H	1.05	2/1709 (0.1%)	1.02	4/2326 (0.2%)
All	All	1.05	3/3397 (0.1%)	1.04	7/4610 (0.2%)

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	L	0	1
2	H	0	1
All	All	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	88	ALA	CA-CB	6.87	1.66	1.52
2	H	12	MET	SD-CE	-5.64	1.46	1.77
1	L	93	GLU	CG-CD	5.51	1.60	1.51

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	57	THR	N-CA-C	6.29	127.97	111.00
2	H	38	LYS	CD-CE-NZ	-6.05	97.78	111.70
1	L	114	SER	N-CA-C	-5.79	95.37	111.00
2	H	178	LEU	CA-CB-CG	5.40	127.71	115.30
1	L	181	LEU	CA-CB-CG	5.33	127.56	115.30
2	H	102	TYR	N-CA-C	-5.28	96.74	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	83	MET	CB-CG-SD	5.07	127.62	112.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	90	TYR	Sidechain
1	L	173	TYR	Sidechain

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	L	1652	0	1604	58	0
2	H	1664	0	1604	103	0
3	L	4	0	0	0	0
4	H	28	0	27	4	0
5	H	56	0	0	4	0
5	L	82	0	0	6	0
All	All	3486	0	3235	152	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 23.

All (152) 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)
2:H:192:GLN:HE21	2:H:192:GLN:HA	1.24	1.00
1:L:80:TYR:O	1:L:83:MET:HG2	1.71	0.90
2:H:192:GLN:NE2	2:H:192:GLN:HA	1.94	0.82
2:H:38:LYS:NZ	2:H:90:TYR:CE2	2.48	0.81
2:H:66:LYS:HB2	2:H:66:LYS:NZ	1.98	0.78
1:L:180:THR:HG23	5:L:518:HOH:O	1.83	0.78
2:H:20:ILE:HG13	2:H:107:THR:HG21	1.65	0.76
2:H:54:SER:O	2:H:56:SER:N	2.18	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:15:LEU:HD12	1:L:16:GLY:N	2.01	0.76
1:L:38:GLN:HE22	2:H:39:GLN:HE22	1.34	0.75
1:L:80:TYR:O	1:L:83:MET:CG	2.36	0.73
2:H:192:GLN:HE21	2:H:192:GLN:CA	1.93	0.73
1:L:187:GLU:HA	1:L:211:ARG:NE	2.04	0.72
2:H:46:GLU:OE2	5:H:416:HOH:O	2.08	0.71
2:H:66:LYS:HB2	2:H:66:LYS:HZ2	1.53	0.71
2:H:10:GLU:HB3	2:H:12:MET:HE1	1.72	0.70
1:L:4:MET:CE	1:L:23:CYS:SG	2.79	0.70
2:H:38:LYS:NZ	2:H:38:LYS:HB2	2.06	0.70
2:H:38:LYS:NZ	2:H:90:TYR:HE2	1.87	0.70
2:H:59:TYR:CD1	2:H:64:LYS:HE2	2.27	0.70
1:L:198:HIS:CD2	1:L:200:GLY:H	2.10	0.69
1:L:4:MET:HE3	1:L:23:CYS:SG	2.33	0.68
2:H:134:GLY:H	2:H:186:SER:HB3	1.58	0.68
1:L:150:VAL:O	1:L:191:VAL:O	2.11	0.68
1:L:1:ASP:HB3	1:L:93:GLU:OE2	1.93	0.68
1:L:145:LYS:NZ	1:L:147:GLN:NE2	2.42	0.68
2:H:101:ASP:CG	4:H:401:HOP:H3	2.15	0.67
2:H:1:GLN:HE21	2:H:2:VAL:HG23	1.60	0.66
1:L:38:GLN:NE2	2:H:39:GLN:HE22	1.93	0.66
2:H:213:PRO:O	2:H:214:LYS:HB2	1.96	0.65
1:L:49:TYR:CZ	2:H:100(B):ASP:HA	2.31	0.65
2:H:129:LYS:HD3	2:H:129:LYS:C	2.17	0.65
2:H:54:SER:C	2:H:56:SER:H	2.00	0.64
2:H:159:LEU:HD21	2:H:182:VAL:HG21	1.80	0.63
2:H:38:LYS:HZ1	2:H:90:TYR:HE2	1.47	0.63
2:H:40:ARG:HD2	2:H:85:GLU:O	1.99	0.63
1:L:186:TYR:O	1:L:211:ARG:HD2	1.99	0.63
2:H:7:SER:OG	2:H:20:ILE:HG12	2.00	0.62
2:H:52:LEU:HB3	2:H:54:SER:O	1.98	0.62
1:L:39:LYS:HE2	1:L:81:GLU:O	1.99	0.62
2:H:129:LYS:NZ	2:H:130:SER:HB2	2.15	0.62
2:H:134:GLY:N	2:H:186:SER:HB3	2.14	0.61
2:H:10:GLU:HB3	2:H:12:MET:CE	2.30	0.61
1:L:193:ALA:HB2	1:L:208:SER:HB3	1.83	0.61
1:L:49:TYR:CD1	2:H:100(A):TYR:CE2	2.89	0.61
2:H:2:VAL:HB	2:H:102:TYR:CE1	2.36	0.61
2:H:32:TYR:CD2	2:H:97:SER:HB2	2.36	0.61
1:L:1:ASP:HB2	1:L:95:PRO:HD2	1.83	0.60
1:L:38:GLN:HE22	2:H:39:GLN:NE2	1.98	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:128:SER:O	2:H:129:LYS:HD3	2.02	0.59
1:L:59:PRO:HG2	5:L:566:HOH:O	2.01	0.59
1:L:145:LYS:HZ1	1:L:147:GLN:NE2	2.00	0.59
2:H:1:GLN:HE21	2:H:2:VAL:CG2	2.16	0.59
2:H:184:VAL:HG11	2:H:194:TYR:CZ	2.38	0.59
2:H:192:GLN:NE2	2:H:192:GLN:CA	2.58	0.58
1:L:192:TYR:O	1:L:208:SER:HB2	2.03	0.58
1:L:198:HIS:HD2	1:L:200:GLY:H	1.51	0.58
1:L:175:LEU:HD23	1:L:175:LEU:C	2.24	0.57
1:L:4:MET:HE2	1:L:23:CYS:SG	2.45	0.57
1:L:49:TYR:HB2	2:H:100(A):TYR:OH	2.05	0.57
2:H:1:GLN:CD	2:H:26:GLY:O	2.43	0.57
2:H:2:VAL:HG13	2:H:27:TYR:HD1	1.69	0.56
2:H:1:GLN:HG2	2:H:26:GLY:HA3	1.88	0.56
2:H:130:SER:OG	2:H:131:THR:N	2.38	0.56
1:L:165:GLU:OE2	5:L:523:HOH:O	2.18	0.55
2:H:2:VAL:HG13	2:H:27:TYR:CD1	2.42	0.54
2:H:17:SER:HB3	2:H:81:GLN:HE21	1.72	0.54
2:H:101:ASP:OD1	4:H:401:HOP:H3	2.08	0.54
2:H:119:PRO:HB3	2:H:145:TYR:HB3	1.90	0.54
2:H:40:ARG:HG3	2:H:88:ALA:HB2	1.90	0.53
1:L:145:LYS:HB3	1:L:197:THR:HB	1.89	0.53
2:H:51:ILE:HA	2:H:56:SER:O	2.08	0.53
2:H:29:PHE:CD2	2:H:76:ASN:HA	2.44	0.52
2:H:38:LYS:HD2	5:H:416:HOH:O	2.10	0.52
2:H:66:LYS:NZ	2:H:66:LYS:CB	2.66	0.52
1:L:95:PRO:O	1:L:97:THR:HG23	2.09	0.52
1:L:193:ALA:CB	1:L:208:SER:HB3	2.40	0.51
1:L:145:LYS:HZ3	1:L:147:GLN:NE2	2.09	0.51
1:L:37:GLN:HB2	1:L:47:LEU:HD11	1.93	0.50
1:L:141:PRO:O	1:L:198:HIS:HE1	1.95	0.50
1:L:165:GLU:HG3	5:L:523:HOH:O	2.10	0.50
1:L:124:GLN:HG2	1:L:129:THR:O	2.12	0.50
2:H:54:SER:C	2:H:56:SER:N	2.63	0.50
1:L:186:TYR:HH	1:L:209:PHE:HE1	1.59	0.49
2:H:126:PRO:HG3	2:H:189:LEU:HD22	1.94	0.49
2:H:99:TYR:CD1	2:H:100:PHE:N	2.80	0.49
2:H:154:TRP:CH2	2:H:196:CYS:HB3	2.48	0.49
2:H:195:ILE:HD12	2:H:195:ILE:N	2.27	0.49
2:H:138:LEU:HD12	2:H:138:LEU:C	2.34	0.48
1:L:161:GLU:CD	5:L:574:HOH:O	2.51	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:192:GLN:CD	2:H:193:THR:H	2.17	0.48
2:H:56:SER:OG	2:H:57:THR:HG23	2.12	0.48
1:L:49:TYR:HD1	2:H:100(A):TYR:CE2	2.32	0.48
2:H:66:LYS:HB2	2:H:66:LYS:HZ3	1.75	0.48
2:H:3:GLN:CG	2:H:25:THR:OG1	2.61	0.48
2:H:186:SER:O	2:H:188:SER:N	2.46	0.48
2:H:61:GLU:HA	2:H:61:GLU:OE1	2.15	0.47
2:H:32:TYR:CE2	2:H:97:SER:HB2	2.49	0.47
2:H:100(A):TYR:CZ	4:H:401:HOP:H18	2.50	0.46
1:L:183:LYS:O	1:L:187:GLU:HG2	2.15	0.46
2:H:66:LYS:HZ2	2:H:66:LYS:CB	2.26	0.46
2:H:154:TRP:HB3	2:H:159:LEU:HD13	1.98	0.46
1:L:161:GLU:HA	1:L:176:SER:O	2.15	0.46
1:L:52:ASN:ND2	5:L:529:HOH:O	2.48	0.46
2:H:146:PHE:HB2	2:H:175:LEU:HD23	1.97	0.45
1:L:150:VAL:HG13	1:L:192:TYR:CE1	2.51	0.45
1:L:124:GLN:O	1:L:127:SER:OG	2.28	0.45
2:H:1:GLN:NE2	2:H:26:GLY:O	2.50	0.45
1:L:38:GLN:OE1	2:H:39:GLN:NE2	2.50	0.45
2:H:144:ASP:HB3	2:H:175:LEU:HD13	1.98	0.45
2:H:192:GLN:NE2	2:H:193:THR:H	2.15	0.45
1:L:185:ASP:HA	1:L:188:LYS:HB2	1.98	0.45
1:L:175:LEU:HD23	1:L:176:SER:N	2.32	0.45
2:H:24:ALA:C	2:H:25:THR:HG23	2.37	0.45
1:L:167:ASP:HB3	1:L:170:ASP:OD1	2.17	0.45
1:L:190:LYS:O	1:L:210:ASN:HA	2.16	0.44
1:L:180:THR:O	1:L:181:LEU:HD13	2.17	0.44
1:L:1:ASP:CB	1:L:93:GLU:OE2	2.65	0.44
2:H:38:LYS:NZ	2:H:38:LYS:CB	2.60	0.44
2:H:20:ILE:HG13	2:H:107:THR:CG2	2.43	0.44
2:H:213:PRO:O	2:H:214:LYS:CB	2.63	0.44
2:H:66:LYS:O	2:H:82:LEU:HA	2.17	0.43
2:H:128:SER:O	2:H:129:LYS:CD	2.65	0.43
2:H:131:THR:O	2:H:133:GLY:N	2.51	0.43
1:L:145:LYS:HZ3	1:L:147:GLN:CD	2.21	0.43
2:H:121:VAL:HG12	2:H:209:LYS:HG3	2.00	0.43
2:H:17:SER:HA	2:H:82:LEU:O	2.19	0.43
2:H:129:LYS:HE2	2:H:129:LYS:O	2.19	0.42
2:H:96:HIS:NE2	5:H:454:HOH:O	2.17	0.42
2:H:143:LYS:HE2	2:H:171:GLN:OE1	2.19	0.42
2:H:7:SER:HG	2:H:20:ILE:HG12	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:6:GLN:HE21	2:H:20:ILE:HD11	1.84	0.42
1:L:23:CYS:HB2	1:L:35:TRP:CH2	2.55	0.42
2:H:19:LYS:HE2	2:H:79:TYR:CD2	2.55	0.42
2:H:38:LYS:HB2	2:H:38:LYS:HZ1	1.81	0.42
2:H:64:LYS:HD2	2:H:64:LYS:HA	1.82	0.42
1:L:58:VAL:HA	1:L:59:PRO:HD3	1.86	0.42
2:H:77:THR:CG2	2:H:79:TYR:CZ	3.03	0.41
2:H:97:SER:HB3	2:H:100:PHE:O	2.20	0.41
1:L:31:SER:C	1:L:50:ARG:NH1	2.74	0.41
2:H:3:GLN:HG2	2:H:25:THR:OG1	2.19	0.41
2:H:148:GLU:HB3	2:H:149:PRO:HA	2.02	0.41
2:H:50:GLU:OE2	5:H:402:HOH:O	2.21	0.41
2:H:20:ILE:HD11	2:H:107:THR:OG1	2.20	0.41
2:H:52:LEU:HA	2:H:52(A):PRO:HD3	1.92	0.41
1:L:54:LEU:HD22	1:L:58:VAL:HB	2.02	0.41
2:H:123:PRO:HD3	2:H:209:LYS:HE3	2.03	0.40
1:L:38:GLN:NE2	2:H:39:GLN:NE2	2.62	0.40
2:H:59:TYR:HB2	2:H:64:LYS:HD3	2.03	0.40
2:H:100(A):TYR:CD2	4:H:401:HOP:H17	2.56	0.40
2:H:77:THR:HG22	2:H:79:TYR:CE1	2.57	0.40
1:L:36:PHE:HE1	1:L:89:LEU:HB3	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	L	209/211 (99%)	202 (97%)	7 (3%)	0	100	100
2	H	219/221 (99%)	195 (89%)	15 (7%)	9 (4%)	3	1
All	All	428/432 (99%)	397 (93%)	22 (5%)	9 (2%)	9	3

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	55	GLY
2	H	56	SER
2	H	62	LYS
2	H	131	THR
2	H	132	SER
2	H	187	SER
2	H	128	SER
2	H	189	LEU
2	H	129	LYS

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	L	189/189 (100%)	172 (91%)	17 (9%)	12	7
2	H	186/186 (100%)	172 (92%)	14 (8%)	17	11
All	All	375/375 (100%)	344 (92%)	31 (8%)	14	8

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	1	ASP
1	L	15	LEU
1	L	30	ASN
1	L	43	SER
1	L	54	LEU
1	L	83	MET
1	L	122	ASP
1	L	123	GLU
1	L	126	LYS
1	L	143	GLU
1	L	149	LYS
1	L	151	ASP
1	L	154	LEU
1	L	165	GLU

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Mol	Chain	Res	Type
1	L	181	LEU
1	L	187	GLU
1	L	210	ASN
2	H	1	GLN
2	H	5	GLN
2	H	20	ILE
2	H	38	LYS
2	H	40	ARG
2	H	43	HIS
2	H	66	LYS
2	H	99	TYR
2	H	101	ASP
2	H	129	LYS
2	H	159	LEU
2	H	172	SER
2	H	184	VAL
2	H	191	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	30	ASN
1	L	38	GLN
1	L	69	GLN
1	L	137	ASN
1	L	147	GLN
1	L	198	HIS
1	L	199	GLN
2	H	1	GLN
2	H	39	GLN
2	H	81	GLN
2	H	192	GLN
2	H	199	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 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	$\# Z > 2$
4	HOP	H	401	-	27,30,30	2.36	11 (40%)	35,40,40	0.80	1 (2%)

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
4	HOP	H	401	-	-	0/17/32/32	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	401	HOP	C2-C1	2.12	1.43	1.38
4	H	401	HOP	C17-C16	2.39	1.43	1.39
4	H	401	HOP	C18-C13	2.79	1.43	1.39
4	H	401	HOP	C15-C16	2.87	1.44	1.39
4	H	401	HOP	C10-C11	3.05	1.57	1.53
4	H	401	HOP	C3-C4	3.12	1.44	1.39
4	H	401	HOP	C12-C7	3.41	1.57	1.53
4	H	401	HOP	C5-C4	3.42	1.44	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	401	HOP	C14-C13	3.50	1.44	1.39
4	H	401	HOP	C13-C10	4.24	1.58	1.51
4	H	401	HOP	C9-C10	4.41	1.59	1.53

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	401	HOP	C23-N22-C20	2.91	128.67	122.15

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	401	HOP	4	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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.