Full wwPDB X-ray Structure Validation Report (i)

Jan 2, 2024 – 04:00 pm GMT

PDB ID	:	4XIX
Title	:	Carbonic anhydrase Cah3 from Chlamydomonas reinhardtii in complex with phosphate.
Authors	:	Hainzl, T.; Grundstrom, C.; Benlloch, R.; Shevela, D.; Shutova, T.; Messinger, J.; Samuelsson, G.; Sauer-Eriksson, A.E.
Deposited on	:	2015-01-08
Resolution	:	2.70 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:

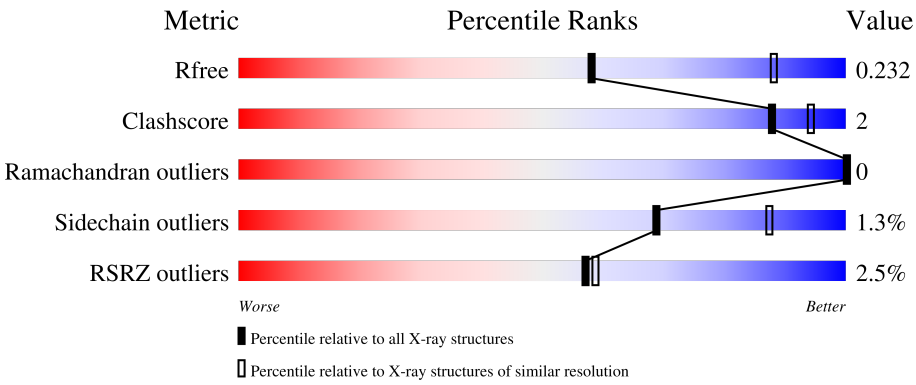
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	239	.% 94% 5%
1	B	239	95% • •
1	C	239	92% 8%
1	D	239	2% 92% 8% •
1	E	239	3% 95% 5%

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Mol	Chain	Length	Quality of chain
1	F	239	8% 95%
1	G	239	3% 94% 5%
1	H	239	3% 93% 7%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 29590 atoms, of which 14584 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 Carbonic anhydrase, alpha type.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	238	3654	1175	1822	311	334	12	0	0	0
1	B	237	3644	1172	1817	310	333	12	0	0	0
1	C	239	3671	1180	1831	312	335	13	0	0	0
1	D	237	3644	1172	1817	310	333	12	0	0	0
1	E	238	3654	1175	1822	311	334	12	0	0	0
1	F	238	3654	1175	1822	311	334	12	0	0	0
1	G	238	3654	1175	1822	311	334	12	0	0	0
1	H	239	3671	1180	1831	312	335	13	0	0	0

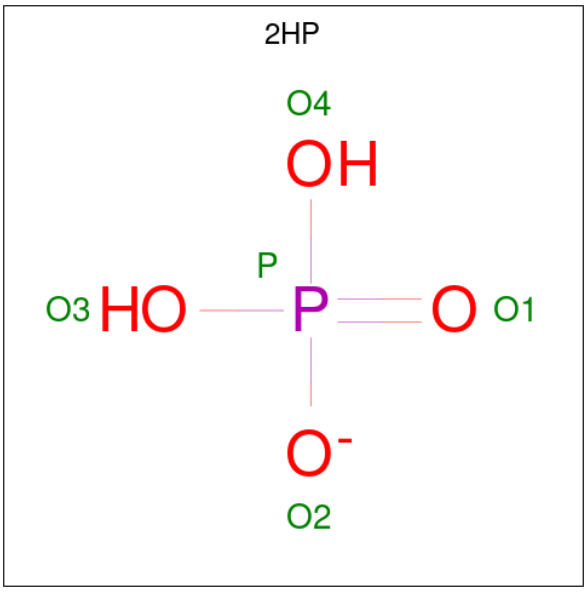
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	72	MET	-	initiating methionine	UNP Q39588
B	72	MET	-	initiating methionine	UNP Q39588
C	72	MET	-	initiating methionine	UNP Q39588
D	72	MET	-	initiating methionine	UNP Q39588
E	72	MET	-	initiating methionine	UNP Q39588
F	72	MET	-	initiating methionine	UNP Q39588
G	72	MET	-	initiating methionine	UNP Q39588
H	72	MET	-	initiating methionine	UNP Q39588

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0
2	E	1	Total Zn 1 1	0	0
2	F	1	Total Zn 1 1	0	0
2	G	1	Total Zn 1 1	0	0
2	H	1	Total Zn 1 1	0	0

- Molecule 3 is DIHYDROGENPHOSPHATE ION (three-letter code: 2HP) (formula: H_2O_4P).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
3	B	1	5	4	1	0	0
			Total	O	P		
3	B	1	5	4	1	0	0
			Total	O	P		
3	B	1	5	4	1	0	0
			Total	O	P		
3	C	1	5	4	1	0	0
			Total	O	P		
3	C	1	5	4	1	0	0
			Total	O	P		
3	C	1	5	4	1	0	0
			Total	O	P		
3	C	1	5	4	1	0	0
			Total	O	P		
3	D	1	5	4	1	0	0
			Total	O	P		
3	D	1	5	4	1	0	0
			Total	O	P		
3	E	1	5	4	1	0	0
			Total	O	P		
3	E	1	5	4	1	0	0
			Total	O	P		
3	E	1	5	4	1	0	0
			Total	O	P		
3	F	1	5	4	1	0	0
			Total	O	P		
3	F	1	5	4	1	0	0
			Total	O	P		
3	G	1	5	4	1	0	0
			Total	O	P		
3	G	1	5	4	1	0	0
			Total	O	P		
3	H	1	5	4	1	0	0
			Total	O	P		
3	H	1	5	4	1	0	0
			Total	O	P		
3	H	1	5	4	1	0	0
			Total	O	P		
3	H	1	5	4	1	0	0

- Molecule 4 is water.

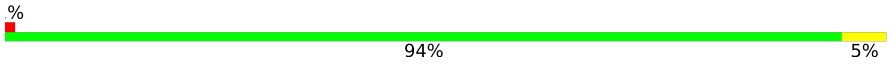
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	46	46	46	0	0
			Total	O		
4	B	34	34	34	0	0
			Total	O		
4	C	36	36	36	0	0
			Total	O		
4	D	16	16	16	0	0
			Total	O		
4	E	21	21	21	0	0
			Total	O		
4	F	15	15	15	0	0
			Total	O		
4	G	20	20	20	0	0
			Total	O		
4	H	28	28	28	0	0

3 Residue-property plots (i)

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

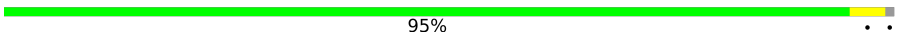
• Molecule 1: Carbonic anhydrase, alpha type

Chain A:



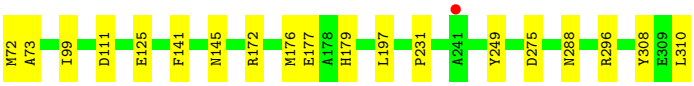
• Molecule 1: Carbonic anhydrase, alpha type

Chain B:



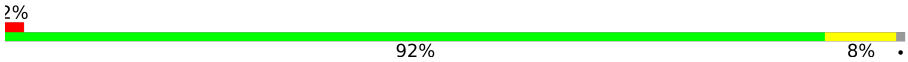
• Molecule 1: Carbonic anhydrase, alpha type

Chain C:



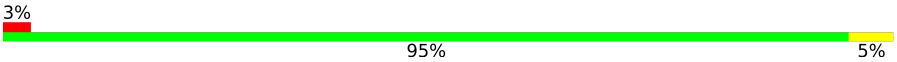
• Molecule 1: Carbonic anhydrase, alpha type

Chain D:



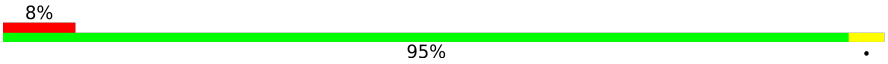
• Molecule 1: Carbonic anhydrase, alpha type

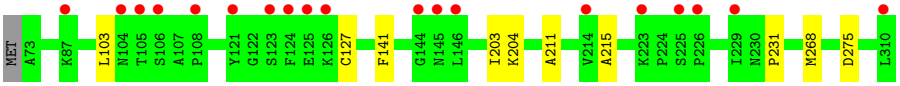
Chain E:



• Molecule 1: Carbonic anhydrase, alpha type

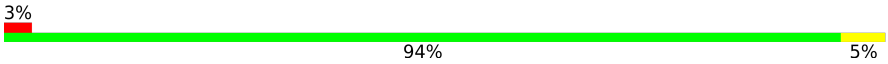
Chain F:





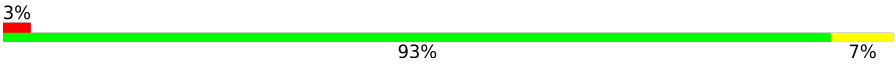
- Molecule 1: Carbonic anhydrase, alpha type

Chain G:



- Molecule 1: Carbonic anhydrase, alpha type

Chain H:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	138.96Å 138.96Å 202.81Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.45 – 2.70 48.45 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.45-2.70) 99.9 (48.45-2.70)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.65 (at 2.69Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.156 , 0.230 0.163 , 0.232	Depositor DCC
R_{free} test set	3178 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	48.5	Xtriage
Anisotropy	0.719	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 40.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.025 for -h,-k,l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	29590	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.88 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.2126e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: 2HP, 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	0.66	0/1884	0.66	0/2561
1	B	0.70	0/1879	0.70	0/2554
1	C	0.69	0/1892	0.73	0/2571
1	D	0.68	0/1879	0.70	0/2554
1	E	0.69	0/1884	0.66	0/2561
1	F	0.62	0/1884	0.67	0/2561
1	G	0.70	0/1884	0.68	0/2561
1	H	0.65	0/1892	0.70	0/2571
All	All	0.67	0/15078	0.69	0/20494

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 (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	1832	1822	1818	7	0
1	B	1827	1817	1813	7	0
1	C	1840	1831	1827	12	0
1	D	1827	1817	1813	9	0
1	E	1832	1822	1818	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	1832	1822	1818	6	0
1	G	1832	1822	1818	8	0
1	H	1840	1831	1827	10	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
3	A	20	0	0	0	0
3	B	15	0	0	0	0
3	C	20	0	0	0	0
3	D	10	0	0	0	0
3	E	15	0	0	0	0
3	F	10	0	0	0	0
3	G	10	0	0	0	0
3	H	20	0	0	0	0
4	A	46	0	0	0	0
4	B	34	0	0	0	0
4	C	36	0	0	0	0
4	D	16	0	0	0	0
4	E	21	0	0	0	0
4	F	15	0	0	0	0
4	G	20	0	0	0	0
4	H	28	0	0	0	0
All	All	15006	14584	14552	59	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 2.

All (59) 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:E:177:GLU:OE2	1:E:179:HIS:NE2	2.28	0.61
1:H:278:ILE:HG22	1:H:282:MET:CE	2.37	0.55
1:B:203:ILE:HB	1:F:203:ILE:HD11	1.90	0.53
1:C:72:MET:HE1	1:E:219:PRO:HB2	1.91	0.53
1:A:177:GLU:OE2	1:A:179:HIS:NE2	2.39	0.53
1:F:103:LEU:HD21	1:F:268:MET:SD	2.48	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:PHE:CE2	1:B:231:PRO:HG3	2.45	0.52
1:C:176:MET:HB2	1:C:197:LEU:HB2	1.93	0.51
1:H:103:LEU:O	1:H:105:THR:N	2.42	0.51
1:A:121:TYR:OH	1:A:182:HIS:NE2	2.40	0.51
1:E:275:ASP:N	1:E:275:ASP:OD1	2.44	0.50
1:C:72:MET:O	1:C:73:ALA:HB3	2.12	0.49
1:H:176:MET:HB2	1:H:197:LEU:HB2	1.94	0.49
1:D:206:PRO:HG2	1:D:233:MET:HG3	1.95	0.48
1:C:177:GLU:OE2	1:C:179:HIS:NE2	2.45	0.48
1:D:247:VAL:HG13	1:D:268:MET:CE	2.44	0.48
1:G:141:PHE:CZ	1:G:231:PRO:HG3	2.48	0.48
1:D:247:VAL:HG13	1:D:268:MET:HE3	1.96	0.48
1:H:130:LEU:HB3	1:H:222:LYS:HG2	1.96	0.48
1:G:148:PHE:CE2	1:G:153:GLU:HG3	2.50	0.47
1:H:275:ASP:N	1:H:275:ASP:OD1	2.47	0.47
1:E:167:HIS:HE1	1:E:264:TRP:CZ3	2.33	0.47
1:H:136:THR:HG21	1:H:160:HIS:HB2	1.96	0.47
1:A:83:PRO:HA	1:A:86:TRP:CD2	2.50	0.46
1:A:275:ASP:OD1	1:A:275:ASP:N	2.48	0.46
1:C:141:PHE:CE1	1:C:231:PRO:HG3	2.50	0.46
1:H:83:PRO:HA	1:H:86:TRP:CD2	2.50	0.46
1:F:203:ILE:HG22	1:F:204:LYS:O	2.15	0.46
1:D:124:PHE:HB2	1:D:229:ILE:O	2.16	0.45
1:B:148:PHE:CE1	1:B:153:GLU:HB2	2.52	0.45
1:F:141:PHE:CZ	1:F:231:PRO:HG3	2.52	0.45
1:D:193:LEU:HD22	1:D:265:PHE:HB2	1.99	0.45
1:B:180:LEU:HD12	1:B:180:LEU:N	2.31	0.44
1:D:83:PRO:HA	1:D:86:TRP:CE2	2.52	0.44
1:E:217:GLU:HG2	1:H:72:MET:HE2	1.99	0.44
1:C:275:ASP:N	1:C:275:ASP:OD1	2.51	0.43
1:B:275:ASP:OD1	1:B:275:ASP:N	2.51	0.43
1:D:83:PRO:HA	1:D:86:TRP:CD2	2.53	0.43
1:A:77:TYR:CE2	1:G:294:ASN:HB2	2.52	0.43
1:G:203:ILE:HG22	1:G:204:LYS:O	2.19	0.43
1:C:172:ARG:HH22	1:H:72:MET:HG2	1.82	0.43
1:G:176:MET:HB2	1:G:197:LEU:HB2	1.99	0.42
1:A:268:MET:HG3	1:A:310:LEU:CD2	2.49	0.42
1:C:99:ILE:HG12	1:C:249:TYR:CZ	2.55	0.42
1:C:125:GLU:OE2	1:C:145:ASN:ND2	2.44	0.42
1:G:230:ASN:HA	1:G:231:PRO:HD3	1.96	0.42
1:A:285:VAL:HG12	1:A:292:ALA:HB3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:275:ASP:OD1	1:G:275:ASP:N	2.53	0.41
1:G:268:MET:HA	1:G:268:MET:HE2	2.02	0.41
1:F:275:ASP:N	1:F:275:ASP:OD1	2.54	0.41
1:B:141:PHE:O	1:C:72:MET:O	2.39	0.41
1:F:211:ALA:O	1:F:215:ALA:HB2	2.21	0.41
1:D:285:VAL:HG12	1:D:292:ALA:HB3	2.03	0.41
1:C:308:TYR:HB3	1:C:310:LEU:HG	2.03	0.41
1:E:220:LEU:HD12	1:E:220:LEU:N	2.36	0.40
1:C:275:ASP:OD2	1:E:276:SER:HA	2.21	0.40
1:B:77:TYR:O	1:H:80:VAL:CG1	2.70	0.40
1:E:83:PRO:HA	1:E:86:TRP:CD2	2.56	0.40
1:D:94:LYS:HB3	1:D:302:ASN:HD22	1.87	0.40

There are no symmetry-related clashes.

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	236/239 (99%)	220 (93%)	16 (7%)	0	100	100
1	B	235/239 (98%)	222 (94%)	13 (6%)	0	100	100
1	C	237/239 (99%)	219 (92%)	18 (8%)	0	100	100
1	D	235/239 (98%)	225 (96%)	10 (4%)	0	100	100
1	E	236/239 (99%)	220 (93%)	16 (7%)	0	100	100
1	F	236/239 (99%)	218 (92%)	18 (8%)	0	100	100
1	G	236/239 (99%)	220 (93%)	16 (7%)	0	100	100
1	H	237/239 (99%)	223 (94%)	14 (6%)	0	100	100
All	All	1888/1912 (99%)	1767 (94%)	121 (6%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	198/199 (100%)	196 (99%)	2 (1%)	76	91
1	B	198/199 (100%)	195 (98%)	3 (2%)	65	86
1	C	199/199 (100%)	196 (98%)	3 (2%)	65	86
1	D	198/199 (100%)	194 (98%)	4 (2%)	55	81
1	E	198/199 (100%)	196 (99%)	2 (1%)	76	91
1	F	198/199 (100%)	197 (100%)	1 (0%)	88	96
1	G	198/199 (100%)	196 (99%)	2 (1%)	76	91
1	H	199/199 (100%)	196 (98%)	3 (2%)	65	86
All	All	1586/1592 (100%)	1566 (99%)	20 (1%)	69	87

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	169	MET
1	A	275	ASP
1	B	130	LEU
1	B	153	GLU
1	B	186	SER
1	C	111	ASP
1	C	288	ASN
1	C	296	ARG
1	D	128	ASP
1	D	158	GLN
1	D	183	LYS
1	D	288	ASN
1	E	275	ASP
1	E	295	THR
1	F	127	CYS
1	G	166	GLU
1	G	275	ASP
1	H	104	ASN
1	H	275	ASP

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Mol	Chain	Res	Type
1	H	303	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	189	ASN
1	E	248	HIS
1	F	138	GLN
1	F	140	ASN
1	G	76	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 32 ligands modelled in this entry, 8 are monoatomic - leaving 24 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	$\#\|Z\| > 2$
3	2HP	C	405	-	4,4,4	1.18	1 (25%)	6,6,6	0.45	0
3	2HP	E	403	-	4,4,4	1.17	1 (25%)	6,6,6	0.44	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	2HP	B	404	-	4,4,4	1.18	1 (25%)	6,6,6	0.45	0
3	2HP	H	405	-	4,4,4	1.17	1 (25%)	6,6,6	0.44	0
3	2HP	C	403	-	4,4,4	1.11	1 (25%)	6,6,6	0.39	0
3	2HP	F	403	-	4,4,4	1.19	1 (25%)	6,6,6	0.45	0
3	2HP	A	402	2	4,4,4	1.17	1 (25%)	6,6,6	0.45	0
3	2HP	G	403	-	4,4,4	1.11	1 (25%)	6,6,6	0.38	0
3	2HP	D	403	-	4,4,4	1.18	1 (25%)	6,6,6	0.44	0
3	2HP	H	403	-	4,4,4	1.19	1 (25%)	6,6,6	0.45	0
3	2HP	G	402	2	4,4,4	1.18	1 (25%)	6,6,6	0.44	0
3	2HP	E	402	2	4,4,4	1.17	1 (25%)	6,6,6	0.44	0
3	2HP	D	402	-	4,4,4	1.19	1 (25%)	6,6,6	0.46	0
3	2HP	B	402	2	4,4,4	1.19	1 (25%)	6,6,6	0.46	0
3	2HP	H	402	2	4,4,4	1.10	1 (25%)	6,6,6	0.38	0
3	2HP	H	404	-	4,4,4	1.18	1 (25%)	6,6,6	0.46	0
3	2HP	A	403	-	4,4,4	1.18	1 (25%)	6,6,6	0.45	0
3	2HP	E	404	-	4,4,4	1.18	1 (25%)	6,6,6	0.45	0
3	2HP	C	404	-	4,4,4	1.18	1 (25%)	6,6,6	0.45	0
3	2HP	B	403	-	4,4,4	1.18	1 (25%)	6,6,6	0.44	0
3	2HP	A	405	-	4,4,4	1.18	1 (25%)	6,6,6	0.44	0
3	2HP	C	402	2	4,4,4	1.16	1 (25%)	6,6,6	0.45	0
3	2HP	F	402	2	4,4,4	1.17	1 (25%)	6,6,6	0.44	0
3	2HP	A	404	-	4,4,4	1.17	1 (25%)	6,6,6	0.44	0

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	2HP	P-O1	2.26	1.56	1.50
3	H	403	2HP	P-O1	2.25	1.56	1.50
3	D	402	2HP	P-O1	2.25	1.56	1.50
3	H	404	2HP	P-O1	2.25	1.56	1.50
3	F	403	2HP	P-O1	2.24	1.56	1.50
3	C	404	2HP	P-O1	2.23	1.56	1.50
3	B	404	2HP	P-O1	2.23	1.56	1.50
3	G	402	2HP	P-O1	2.22	1.56	1.50
3	A	403	2HP	P-O1	2.22	1.56	1.50
3	E	404	2HP	P-O1	2.22	1.56	1.50
3	A	405	2HP	P-O1	2.22	1.56	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	402	2HP	P-O1	2.22	1.56	1.50
3	B	403	2HP	P-O1	2.21	1.56	1.50
3	D	403	2HP	P-O1	2.21	1.56	1.50
3	C	405	2HP	P-O1	2.21	1.56	1.50
3	E	403	2HP	P-O1	2.21	1.56	1.50
3	A	402	2HP	P-O1	2.21	1.56	1.50
3	H	405	2HP	P-O1	2.20	1.56	1.50
3	A	404	2HP	P-O1	2.20	1.56	1.50
3	F	402	2HP	P-O1	2.19	1.56	1.50
3	C	402	2HP	P-O1	2.18	1.56	1.50
3	C	403	2HP	P-O1	2.05	1.55	1.50
3	G	403	2HP	P-O1	2.03	1.55	1.50
3	H	402	2HP	P-O1	2.02	1.55	1.50

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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)

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	238/239 (99%)	-0.15	3 (1%)	77	78	12, 27, 47, 63	0
1	B	237/239 (99%)	-0.19	1 (0%)	92	93	10, 25, 43, 66	0
1	C	239/239 (100%)	-0.18	1 (0%)	92	93	13, 24, 43, 54	0
1	D	237/239 (99%)	-0.06	4 (1%)	70	72	14, 30, 51, 60	0
1	E	238/239 (99%)	0.03	7 (2%)	51	52	13, 28, 53, 83	0
1	F	238/239 (99%)	0.18	19 (7%)	12	10	13, 37, 64, 87	0
1	G	238/239 (99%)	-0.05	6 (2%)	57	59	10, 29, 54, 73	0
1	H	239/239 (100%)	-0.03	7 (2%)	51	52	13, 28, 55, 73	0
All	All	1904/1912 (99%)	-0.06	48 (2%)	57	59	10, 28, 54, 87	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	310	LEU	5.2
1	D	310	LEU	5.2
1	F	104	ASN	4.7
1	F	310	LEU	4.7
1	E	106	SER	4.6
1	D	92	THR	3.9
1	F	226	PRO	3.7
1	A	310	LEU	3.6
1	G	126	LYS	3.4
1	F	105	THR	3.3
1	F	229	ILE	3.2
1	A	92	THR	3.2
1	H	104	ASN	3.0
1	F	125	GLU	3.0
1	G	123	SER	2.9
1	F	144	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	G	310	LEU	2.8
1	D	95	ARG	2.8
1	F	123	SER	2.8
1	E	103	LEU	2.7
1	H	310	LEU	2.6
1	F	225	SER	2.5
1	C	241	ALA	2.4
1	F	87	LYS	2.4
1	F	126	LYS	2.4
1	H	122	GLY	2.4
1	A	73	ALA	2.4
1	H	103	LEU	2.3
1	E	226	PRO	2.3
1	B	310	LEU	2.3
1	F	223	LYS	2.3
1	F	108	PRO	2.3
1	F	106	SER	2.2
1	E	125	GLU	2.2
1	E	104	ASN	2.2
1	F	124	PHE	2.2
1	G	105	THR	2.2
1	H	226	PRO	2.1
1	H	106	SER	2.1
1	G	229	ILE	2.1
1	G	104	ASN	2.1
1	D	89	VAL	2.1
1	F	121	TYR	2.1
1	F	214	VAL	2.1
1	F	145	ASN	2.0
1	E	105	THR	2.0
1	F	146	LEU	2.0
1	H	126	LYS	2.0

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 monosaccharides 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. 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	B-factors(Å^2)	Q<0.9
3	2HP	D	403	5/5	0.88	0.24	35,47,78,89	0
3	2HP	C	405	5/5	0.89	0.18	37,44,58,72	0
3	2HP	B	403	5/5	0.90	0.25	15,21,28,44	5
3	2HP	E	404	5/5	0.90	0.23	13,22,28,28	5
3	2HP	B	404	5/5	0.91	0.32	21,32,36,38	5
3	2HP	A	403	5/5	0.91	0.28	38,46,52,63	0
3	2HP	E	403	5/5	0.94	0.17	44,55,63,68	0
3	2HP	F	403	5/5	0.94	0.14	41,44,52,66	0
3	2HP	H	405	5/5	0.95	0.08	41,41,47,66	0
3	2HP	A	405	5/5	0.96	0.16	29,32,36,41	0
3	2HP	G	403	5/5	0.96	0.13	35,44,51,64	0
3	2HP	H	402	5/5	0.96	0.19	25,27,35,39	0
3	2HP	H	404	5/5	0.96	0.20	25,54,57,58	0
3	2HP	B	402	5/5	0.96	0.29	21,25,40,41	0
3	2HP	E	402	5/5	0.97	0.26	10,23,37,41	0
3	2HP	A	402	5/5	0.97	0.19	18,21,31,36	0
3	2HP	G	402	5/5	0.98	0.19	20,22,26,32	0
3	2HP	C	402	5/5	0.98	0.21	25,26,37,43	0
3	2HP	C	403	5/5	0.98	0.13	20,23,25,30	0
3	2HP	H	403	5/5	0.98	0.15	29,29,30,33	0
3	2HP	F	402	5/5	0.98	0.20	22,23,36,36	0
3	2HP	C	404	5/5	0.98	0.14	15,16,20,24	0
2	ZN	C	401	1/1	0.99	0.15	24,24,24,24	0
3	2HP	A	404	5/5	0.99	0.15	28,29,34,38	0
3	2HP	D	402	5/5	0.99	0.19	21,25,32,35	0
2	ZN	H	401	1/1	0.99	0.16	19,19,19,19	0
2	ZN	B	401	1/1	0.99	0.17	20,20,20,20	0
2	ZN	A	401	1/1	1.00	0.15	18,18,18,18	0
2	ZN	D	401	1/1	1.00	0.15	26,26,26,26	0
2	ZN	E	401	1/1	1.00	0.16	19,19,19,19	0
2	ZN	F	401	1/1	1.00	0.16	26,26,26,26	0
2	ZN	G	401	1/1	1.00	0.17	16,16,16,16	0

6.5 Other polymers (i)

There are no such residues in this entry.