Full wwPDB X-ray Structure Validation Report (i)

Oct 17, 2021 – 01:13 AM EDT

PDB ID	:	1QW9
Title	:	Crystal structure of a family 51 alpha-L-arabinofuranosidase in complex with 4-nitrophenyl-Ara
Authors	:	Hoevel, K.; Shallom, D.; Niefind, K.; Belakhov, V.; Shoham, G.; Bassov, T.; Shoham, Y.; Schomburg, D.
Deposited on	:	2003-09-01
Resolution	:	1.20 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 following versions of software and data (see references (i)) were used in the production of this report:

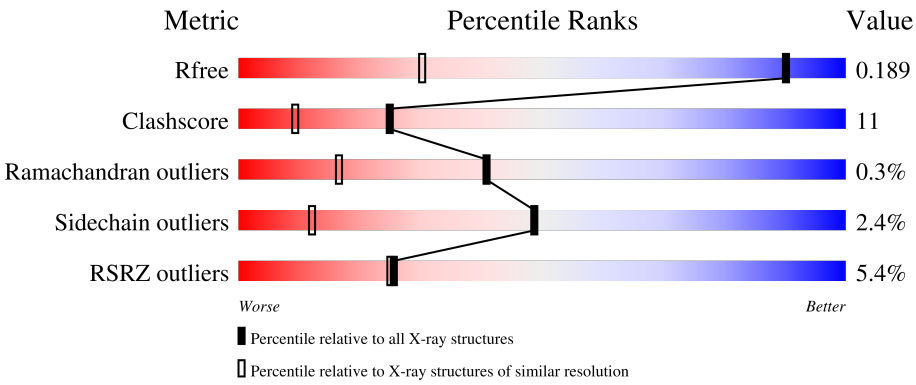
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

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 1.20 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	1223 (1.22-1.18)
Clashscore	141614	1286 (1.22-1.18)
Ramachandran outliers	138981	1240 (1.22-1.18)
Sidechain outliers	138945	1239 (1.22-1.18)
RSRZ outliers	127900	1200 (1.22-1.18)

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	502	6% 81% 16% • •
1	B	502	4% 84% 14% • •

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 9598 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.

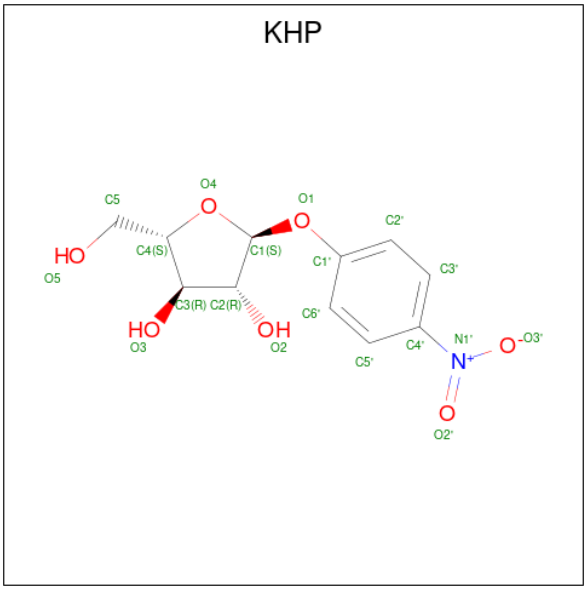
- Molecule 1 is a protein called Alpha-L-arabinofuranosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	497	3986	2540	680	746	20	0	0	0
1	B	497	3986	2540	680	746	20	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	175	ALA	GLU	engineered mutation	UNP Q9XBQ3
B	175	ALA	GLU	engineered mutation	UNP Q9XBQ3

- Molecule 2 is 4-nitrophenyl alpha-L-arabinofuranoside (three-letter code: KHP) (formula: $C_{11}H_{13}NO_7$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	19	11	1	7	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	B	1	19	11	1	7	0	0

- Molecule 3 is water.

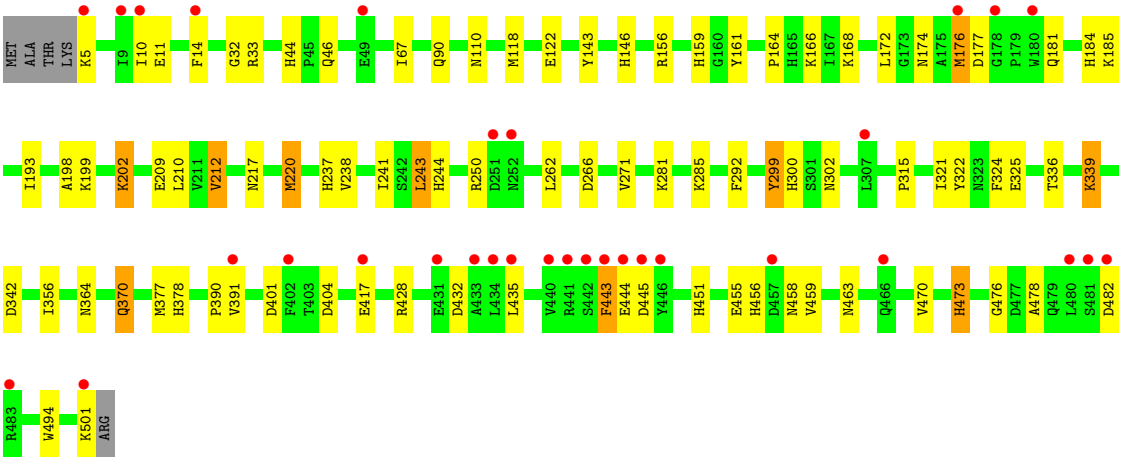
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	805	805	805	0	0
			Total	O		
3	B	783	783	783	0	0

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.

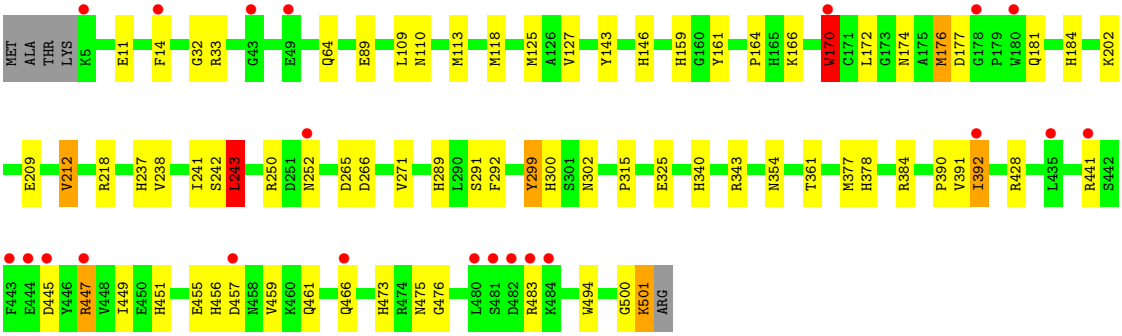
• Molecule 1: Alpha-L-arabinofuranosidase

Chain A:



• Molecule 1: Alpha-L-arabinofuranosidase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α, β, γ	179.31Å 179.31Å 100.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 1.20 44.83 – 1.20	Depositor EDS
% Data completeness (in resolution range)	98.8 (50.00-1.20) 98.8 (44.83-1.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.65 (at 1.20Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.160 , 0.179 0.172 , 0.189	Depositor DCC
R_{free} test set	18644 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å^2)	13.6	Xtriage
Anisotropy	0.055	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 40.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.008 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	9598	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.93% of the height of the origin peak. No significant pseudotranslation is detected.*

[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: KHP

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	0.65	3/4087 (0.1%)	0.85	14/5553 (0.3%)
1	B	0.66	6/4087 (0.1%)	0.88	15/5553 (0.3%)
All	All	0.65	9/8174 (0.1%)	0.87	29/11106 (0.3%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	170	TRP	CB-CG	-10.08	1.32	1.50
1	B	176	MET	CB-CG	7.59	1.75	1.51
1	A	220	MET	SD-CE	-6.75	1.40	1.77
1	A	202	LYS	CD-CE	6.66	1.68	1.51
1	A	212	VAL	CB-CG1	-6.64	1.39	1.52
1	B	170	TRP	CG-CD2	6.57	1.54	1.43
1	B	212	VAL	CB-CG1	-6.31	1.39	1.52
1	B	209	GLU	CD-OE1	-6.09	1.19	1.25
1	B	89	GLU	CD-OE1	5.38	1.31	1.25

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	243	LEU	CA-CB-CG	10.63	139.74	115.30
1	B	177	ASP	CB-CG-OD2	9.56	126.90	118.30
1	A	33	ARG	NE-CZ-NH1	8.23	124.42	120.30
1	B	33	ARG	NE-CZ-NH2	-8.22	116.19	120.30
1	B	170	TRP	CB-CG-CD1	-7.95	116.67	127.00
1	B	176	MET	CG-SD-CE	-7.72	87.84	100.20
1	B	170	TRP	CB-CG-CD2	7.69	136.60	126.60
1	B	170	TRP	CG-CD2-CE3	7.44	140.59	133.90
1	B	33	ARG	NE-CZ-NH1	7.42	124.01	120.30
1	B	170	TRP	CE2-CD2-CG	-6.87	101.81	107.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	457	ASP	CB-CG-OD2	6.67	124.30	118.30
1	B	384	ARG	NE-CZ-NH1	6.59	123.60	120.30
1	B	176	MET	CA-CB-CG	6.46	124.29	113.30
1	A	177	ASP	CB-CG-OD1	6.37	124.03	118.30
1	A	281	LYS	CD-CE-NZ	-6.19	97.45	111.70
1	B	428	ARG	NE-CZ-NH1	5.99	123.29	120.30
1	B	428	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	B	265	ASP	CB-CG-OD2	5.93	123.64	118.30
1	A	33	ARG	NE-CZ-NH2	-5.90	117.35	120.30
1	A	445	ASP	CB-CG-OD2	5.76	123.48	118.30
1	A	342	ASP	CB-CG-OD1	5.68	123.41	118.30
1	A	482	ASP	CB-CG-OD2	5.50	123.25	118.30
1	A	432	ASP	CB-CG-OD2	5.36	123.12	118.30
1	A	243	LEU	CB-CA-C	-5.23	100.26	110.20
1	A	404	ASP	CB-CG-OD2	5.23	123.00	118.30
1	A	428	ARG	NE-CZ-NH2	-5.18	117.71	120.30
1	A	266	ASP	CB-CG-OD2	5.16	122.94	118.30
1	A	156	ARG	NE-CZ-NH1	5.04	122.82	120.30
1	A	401	ASP	CB-CG-OD2	5.02	122.82	118.30

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	3986	0	3890	90	0
1	B	3986	0	3890	84	0
2	A	19	0	4	0	0
2	B	19	0	4	0	0
3	A	805	0	0	38	0
3	B	783	0	0	36	1
All	All	9598	0	7788	170	1

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 11.

All (170) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:176:MET:CB	1:B:176:MET:CG	1.75	1.64
1:A:377:MET:HG2	3:A:2482:HOH:O	1.24	1.36
1:B:377:MET:HG2	3:B:2459:HOH:O	1.25	1.29
1:A:243:LEU:HD13	3:A:2487:HOH:O	1.28	1.29
1:A:118:MET:HG3	3:A:2484:HOH:O	1.19	1.28
1:B:166:LYS:HG2	3:B:2351:HOH:O	1.38	1.20
1:A:176:MET:CE	3:A:2486:HOH:O	1.85	1.19
1:B:501:LYS:HA	3:B:2350:HOH:O	1.41	1.17
1:B:377:MET:CG	3:B:2459:HOH:O	1.83	1.15
1:A:118:MET:SD	3:A:2484:HOH:O	2.03	1.14
1:A:176:MET:SD	3:A:2486:HOH:O	2.05	1.14
1:A:176:MET:HG3	3:A:2283:HOH:O	1.48	1.10
1:A:377:MET:CG	3:A:2482:HOH:O	1.84	1.05
1:A:118:MET:CG	3:A:2484:HOH:O	1.78	1.04
1:A:377:MET:SD	3:A:2482:HOH:O	2.16	1.01
1:B:170:TRP:CE3	3:B:2458:HOH:O	2.13	1.01
1:B:377:MET:SD	3:B:2459:HOH:O	2.17	0.98
1:A:166:LYS:HG2	3:A:2290:HOH:O	1.67	0.94
1:A:14:PHE:CZ	1:B:391:VAL:HG21	2.01	0.94
1:A:176:MET:HE3	3:A:2486:HOH:O	1.55	0.92
1:B:170:TRP:HB2	3:B:2458:HOH:O	1.68	0.92
1:B:473:HIS:HD2	1:B:475:ASN:H	1.20	0.88
1:B:202:LYS:HD2	3:B:2208:HOH:O	1.75	0.87
1:A:377:MET:CE	3:A:2482:HOH:O	2.23	0.86
1:A:176:MET:CG	3:A:2283:HOH:O	2.11	0.85
1:B:377:MET:CE	3:B:2459:HOH:O	2.26	0.84
1:B:176:MET:CB	1:B:176:MET:SD	2.67	0.83
1:B:170:TRP:CD2	3:B:2458:HOH:O	2.31	0.82
1:A:14:PHE:CE2	1:B:391:VAL:HG21	2.15	0.81
1:A:176:MET:CG	3:A:2486:HOH:O	2.25	0.80
1:B:392:ILE:HD11	3:B:2039:HOH:O	1.79	0.80
1:A:212:VAL:CG1	1:A:238:VAL:HG11	2.12	0.79
1:A:174:ASN:O	1:A:176:MET:HG2	1.82	0.79
1:B:392:ILE:HG21	3:B:2465:HOH:O	1.83	0.79
1:A:243:LEU:HD11	1:A:271:VAL:HG11	1.65	0.79
1:B:174:ASN:O	1:B:176:MET:HG3	1.82	0.78
1:B:212:VAL:CG1	1:B:238:VAL:HG11	2.13	0.78
1:B:354:ASN:HD21	1:B:361:THR:H	1.29	0.78
1:A:202:LYS:HE2	1:A:237:HIS:O	1.84	0.77
1:A:378:HIS:HD2	1:A:494:TRP:HE1	1.32	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:501:LYS:CA	3:B:2350:HOH:O	2.09	0.77
1:B:377:MET:HE2	3:B:2459:HOH:O	1.83	0.76
1:B:378:HIS:HD2	1:B:494:TRP:HE1	1.32	0.75
1:A:244:HIS:CE1	3:A:1931:HOH:O	2.41	0.74
1:B:501:LYS:CB	3:B:2350:HOH:O	2.37	0.72
1:A:463:ASN:HD21	1:A:470:VAL:H	1.40	0.70
1:B:143:TYR:OH	1:B:159:HIS:HD2	1.75	0.69
1:B:390:PRO:HA	3:B:2462:HOH:O	1.92	0.69
1:B:127:VAL:HG13	1:B:170:TRP:HE1	1.58	0.69
1:B:456:HIS:CE1	1:B:461:GLN:HG2	2.28	0.69
1:B:170:TRP:CG	3:B:2458:HOH:O	2.45	0.68
1:B:250:ARG:HH21	1:B:302:ASN:HD21	1.41	0.68
1:A:364:ASN:CG	3:A:2136:HOH:O	2.33	0.67
1:A:443:PHE:CE2	3:A:2231:HOH:O	2.48	0.67
1:B:449:ILE:HD12	3:B:2437:HOH:O	1.95	0.67
1:A:243:LEU:HB2	3:A:2487:HOH:O	1.94	0.66
1:A:250:ARG:HH21	1:A:302:ASN:HD21	1.41	0.66
1:A:443:PHE:CD2	3:A:2231:HOH:O	2.48	0.66
1:B:266:ASP:OD2	3:B:2344:HOH:O	2.15	0.65
1:A:90:GLN:NE2	3:A:2221:HOH:O	2.30	0.65
1:A:185:LYS:HZ2	1:A:193:ILE:HD13	1.62	0.65
1:B:170:TRP:HE3	3:B:2458:HOH:O	1.61	0.64
1:A:377:MET:HE2	3:A:2482:HOH:O	1.88	0.64
1:B:127:VAL:HG13	1:B:170:TRP:NE1	2.14	0.63
1:A:299:TYR:CZ	1:A:300:HIS:CE1	2.86	0.63
1:B:445:ASP:O	1:B:501:LYS:HB3	1.99	0.63
1:A:243:LEU:CG	3:A:2487:HOH:O	2.44	0.62
1:B:170:TRP:CB	3:B:2458:HOH:O	2.36	0.62
1:B:202:LYS:HD3	3:B:2183:HOH:O	1.99	0.61
1:B:456:HIS:HD2	3:B:2189:HOH:O	1.83	0.61
1:A:11:GLU:CG	1:A:14:PHE:HD1	2.14	0.61
1:A:14:PHE:CE2	1:B:391:VAL:CG2	2.82	0.61
1:A:164:PRO:HG2	1:A:166:LYS:HE2	1.84	0.59
1:A:11:GLU:HG2	1:A:14:PHE:HD1	1.69	0.58
1:A:378:HIS:CD2	1:A:494:TRP:HE1	2.18	0.58
1:A:237:HIS:HE1	3:A:2249:HOH:O	1.86	0.58
1:B:378:HIS:CD2	1:B:494:TRP:HE1	2.19	0.58
1:B:11:GLU:HG2	1:B:14:PHE:HD1	1.69	0.58
1:A:299:TYR:CE1	1:A:300:HIS:CE1	2.92	0.57
1:B:500:GLY:O	1:B:501:LYS:HB2	2.03	0.57
1:A:456:HIS:HD2	1:A:458:ASN:H	1.53	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:ASN:HD22	1:A:181:GLN:HE22	1.54	0.56
1:A:143:TYR:OH	1:A:159:HIS:HD2	1.89	0.56
1:B:176:MET:O	1:B:184:HIS:HD2	1.88	0.56
1:A:10:ILE:HG22	1:A:443:PHE:HZ	1.70	0.55
1:B:146:HIS:HD2	3:B:1858:HOH:O	1.90	0.55
1:B:451:HIS:CG	1:B:476:GLY:HA3	2.42	0.55
1:B:174:ASN:HD22	1:B:181:GLN:HE22	1.54	0.55
1:B:237:HIS:HE1	3:B:2236:HOH:O	1.89	0.54
1:B:212:VAL:HG11	1:B:238:VAL:HG11	1.88	0.54
1:B:455:GLU:OE2	1:B:473:HIS:HE1	1.91	0.54
1:B:237:HIS:HD2	3:B:2133:HOH:O	1.91	0.53
1:A:11:GLU:HG2	1:A:14:PHE:CD1	2.43	0.53
1:B:172:LEU:HD12	1:B:212:VAL:HG12	1.90	0.53
1:A:198:ALA:O	1:A:202:LYS:HE3	2.09	0.53
1:B:500:GLY:O	1:B:501:LYS:CB	2.57	0.53
1:A:159:HIS:HE1	3:A:1822:HOH:O	1.91	0.53
1:B:11:GLU:CG	1:B:14:PHE:HD1	2.22	0.53
1:B:164:PRO:HG2	1:B:166:LYS:HE2	1.91	0.53
1:A:237:HIS:HD2	3:A:2532:HOH:O	1.91	0.52
1:A:185:LYS:NZ	1:A:193:ILE:HD13	2.24	0.52
1:A:44:HIS:CD2	1:A:46:GLN:H	2.26	0.52
1:A:391:VAL:N	3:A:2318:HOH:O	2.41	0.52
1:B:11:GLU:HG2	1:B:14:PHE:CD1	2.45	0.52
1:B:456:HIS:HE1	1:B:461:GLN:HG2	1.71	0.51
1:A:455:GLU:OE2	1:A:473:HIS:HE1	1.93	0.51
1:A:217:ASN:O	1:A:220:MET:HG2	2.10	0.51
1:B:456:HIS:CD2	3:B:2189:HOH:O	2.60	0.50
1:A:198:ALA:HB1	1:A:210:LEU:CD1	2.41	0.50
1:A:262:LEU:HD23	3:A:1954:HOH:O	2.12	0.50
1:B:146:HIS:HE1	3:B:1939:HOH:O	1.93	0.50
1:A:336:THR:HA	1:A:339:LYS:HG2	1.93	0.50
1:A:322:TYR:H	1:A:370:GLN:NE2	2.10	0.50
1:A:451:HIS:HD1	1:A:478:ALA:H	1.60	0.49
1:B:441:ARG:HG2	1:B:483:ARG:HD3	1.94	0.49
1:A:185:LYS:NZ	1:A:193:ILE:CD1	2.77	0.48
1:A:146:HIS:HE1	3:A:1908:HOH:O	1.96	0.48
1:B:451:HIS:CD2	1:B:476:GLY:HA3	2.48	0.48
1:A:199:LYS:NZ	3:A:1759:HOH:O	2.41	0.48
1:A:322:TYR:H	1:A:370:GLN:HE22	1.62	0.47
1:A:10:ILE:HG22	1:A:443:PHE:CZ	2.48	0.47
1:A:168:LYS:HE3	1:A:209:GLU:HG3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:GLN:NE2	3:B:2424:HOH:O	2.45	0.47
1:B:125:MET:O	1:B:170:TRP:CD1	2.68	0.47
1:A:198:ALA:CB	1:A:210:LEU:HD13	2.45	0.47
1:A:243:LEU:HD22	3:A:2487:HOH:O	2.14	0.46
1:B:243:LEU:HG	1:B:271:VAL:HG21	1.97	0.46
1:A:176:MET:HG2	3:A:2486:HOH:O	2.00	0.46
1:A:378:HIS:HE1	3:A:1979:HOH:O	1.98	0.46
1:B:473:HIS:CD2	1:B:475:ASN:H	2.13	0.46
1:B:354:ASN:ND2	1:B:361:THR:H	2.04	0.45
1:B:184:HIS:HE1	3:B:2107:HOH:O	1.98	0.45
1:B:325:GLU:HB3	1:B:459:VAL:HB	1.98	0.45
1:B:340:HIS:HD2	3:B:1801:HOH:O	1.98	0.44
1:A:172:LEU:HD12	1:A:212:VAL:HG12	1.98	0.44
1:A:321:ILE:HA	1:A:370:GLN:HE22	1.83	0.44
1:B:32:GLY:HA2	1:B:315:PRO:O	2.18	0.44
1:B:118:MET:HA	1:B:118:MET:CE	2.47	0.44
1:A:166:LYS:NZ	3:A:2534:HOH:O	2.13	0.44
1:A:435:LEU:HD22	1:A:435:LEU:N	2.32	0.44
1:B:289:HIS:HD2	1:B:343:ARG:HE	1.65	0.44
1:A:456:HIS:CD2	1:A:458:ASN:H	2.33	0.44
1:B:110:ASN:HB3	1:B:161:TYR:CZ	2.52	0.44
1:A:176:MET:O	1:A:184:HIS:HD2	2.02	0.43
1:A:146:HIS:HD2	3:A:1784:HOH:O	2.01	0.43
1:A:325:GLU:HB3	1:A:459:VAL:HB	2.00	0.43
1:A:390:PRO:O	3:A:2513:HOH:O	2.21	0.43
1:A:184:HIS:HE1	3:A:1924:HOH:O	2.01	0.43
1:A:243:LEU:CD1	3:A:2487:HOH:O	2.09	0.43
1:A:110:ASN:HB3	1:A:161:TYR:CZ	2.54	0.43
1:A:324:PHE:CZ	1:A:455:GLU:HA	2.54	0.42
1:B:176:MET:O	1:B:184:HIS:CD2	2.71	0.42
1:A:463:ASN:ND2	1:A:470:VAL:H	2.13	0.42
1:A:32:GLY:HA2	1:A:315:PRO:O	2.19	0.42
1:B:143:TYR:OH	1:B:159:HIS:CD2	2.65	0.42
1:B:501:LYS:CG	3:B:2350:HOH:O	2.65	0.42
1:A:185:LYS:NZ	1:A:193:ILE:HG21	2.35	0.42
1:B:166:LYS:NZ	3:B:2052:HOH:O	2.48	0.42
1:A:451:HIS:CG	1:A:476:GLY:HA3	2.55	0.42
1:B:159:HIS:HE1	3:B:1859:HOH:O	2.03	0.41
1:B:447:ARG:HE	1:B:501:LYS:HD3	1.86	0.41
1:A:212:VAL:HG13	1:A:238:VAL:HG11	1.95	0.41
1:A:473:HIS:CD2	1:A:473:HIS:C	2.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:299:TYR:CZ	1:B:300:HIS:CE1	3.09	0.40
1:B:109:LEU:O	1:B:113:MET:HG2	2.21	0.40
1:A:14:PHE:CZ	1:B:391:VAL:CG2	2.89	0.40
1:A:67:ILE:HG13	1:A:122:GLU:HG3	2.03	0.40
1:B:64:GLN:NE2	3:B:1847:HOH:O	2.54	0.40
1:B:218:ARG:HG2	3:B:2019:HOH:O	2.22	0.40
1:B:242:SER:HA	1:B:291:SER:O	2.21	0.40

All (1) 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)
3:B:1871:HOH:O	3:B:2530:HOH:O[2_655]	1.77	0.43

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	495/502 (99%)	478 (97%)	15 (3%)	2 (0%)	34	11
1	B	495/502 (99%)	479 (97%)	15 (3%)	1 (0%)	47	19
All	All	990/1004 (99%)	957 (97%)	30 (3%)	3 (0%)	41	16

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	299	TYR
1	A	299	TYR
1	A	356	ILE

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	429/433 (99%)	417 (97%)	12 (3%)	43 8
1	B	429/433 (99%)	420 (98%)	9 (2%)	53 15
All	All	858/866 (99%)	837 (98%)	21 (2%)	49 12

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	176	MET
1	A	241	ILE
1	A	285	LYS
1	A	292	PHE
1	A	339	LYS
1	A	370	GLN
1	A	417	GLU
1	A	443	PHE
1	A	444	GLU
1	A	473	HIS
1	A	501	LYS
1	B	170	TRP
1	B	241	ILE
1	B	243	LEU
1	B	252	ASN
1	B	292	PHE
1	B	392	ILE
1	B	447	ARG
1	B	466	GLN
1	B	501	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	HIS

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Mol	Chain	Res	Type
1	A	64	GLN
1	A	146	HIS
1	A	159	HIS
1	A	181	GLN
1	A	184	HIS
1	A	237	HIS
1	A	244	HIS
1	A	245	GLN
1	A	302	ASN
1	A	370	GLN
1	A	378	HIS
1	A	456	HIS
1	A	463	ASN
1	A	473	HIS
1	A	495	ASN
1	B	46	GLN
1	B	64	GLN
1	B	146	HIS
1	B	159	HIS
1	B	181	GLN
1	B	184	HIS
1	B	237	HIS
1	B	252	ASN
1	B	289	HIS
1	B	302	ASN
1	B	340	HIS
1	B	354	ASN
1	B	378	HIS
1	B	456	HIS
1	B	473	HIS

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)

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	KHP	B	1751	-	19,20,20	1.95	6 (31%)	26,28,28	0.91	1 (3%)
2	KHP	A	1750	-	19,20,20	1.98	7 (36%)	26,28,28	1.01	2 (7%)

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	KHP	B	1751	-	-	2/7/26/26	0/2/2/2
2	KHP	A	1750	-	-	0/7/26/26	0/2/2/2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1751	KHP	C4'-N1'	-4.58	1.34	1.45
2	A	1750	KHP	C4'-N1'	-4.40	1.34	1.45
2	A	1750	KHP	O5-C5	-3.23	1.28	1.42
2	A	1750	KHP	O1-C1'	3.21	1.44	1.38
2	B	1751	KHP	O1-C1'	2.92	1.44	1.38
2	B	1751	KHP	O5-C5	-2.91	1.30	1.42
2	A	1750	KHP	O1-C1	2.67	1.45	1.41
2	B	1751	KHP	C3'-C4'	-2.47	1.34	1.38
2	B	1751	KHP	C3'-C2'	-2.29	1.34	1.38
2	B	1751	KHP	C6'-C5'	-2.28	1.34	1.38
2	A	1750	KHP	C3'-C4'	-2.22	1.34	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1750	KHP	C3'-C2'	-2.18	1.34	1.38
2	A	1750	KHP	C5'-C4'	-2.07	1.35	1.38

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1750	KHP	C1-C2-C3	2.82	105.88	102.30
2	B	1751	KHP	O5-C5-C4	2.27	119.06	111.29
2	A	1750	KHP	O5-C5-C4	2.16	118.69	111.29

There are no chirality outliers.

All (2) torsion outliers are listed below:

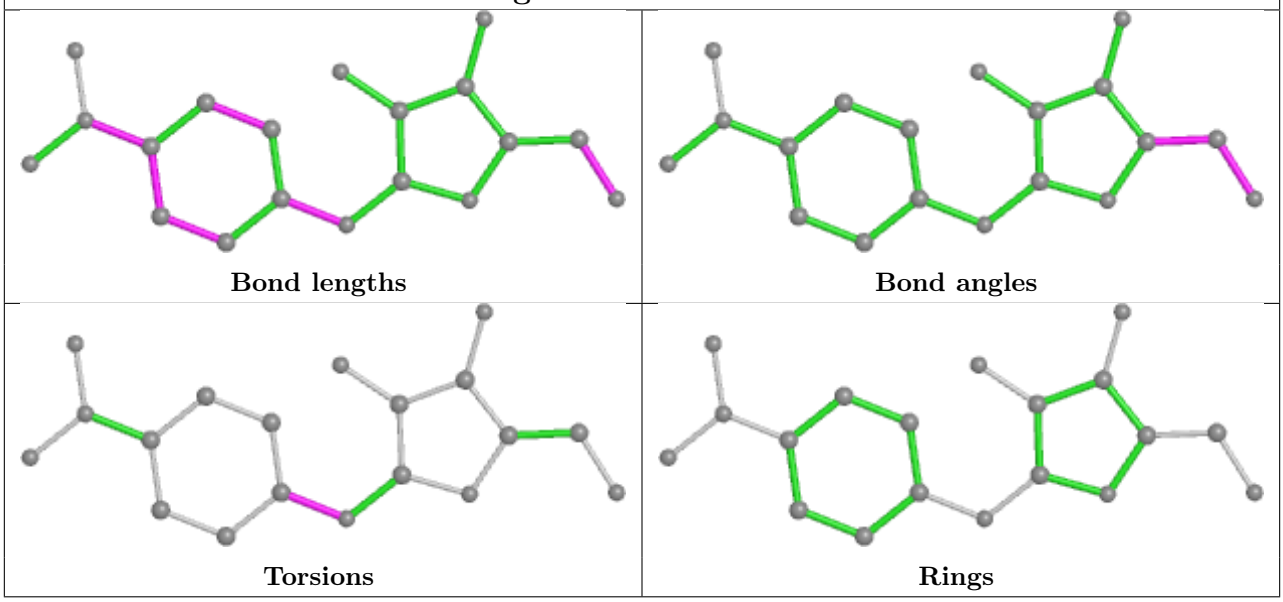
Mol	Chain	Res	Type	Atoms
2	B	1751	KHP	C2'-C1'-O1-C1
2	B	1751	KHP	C6'-C1'-O1-C1

There are no ring outliers.

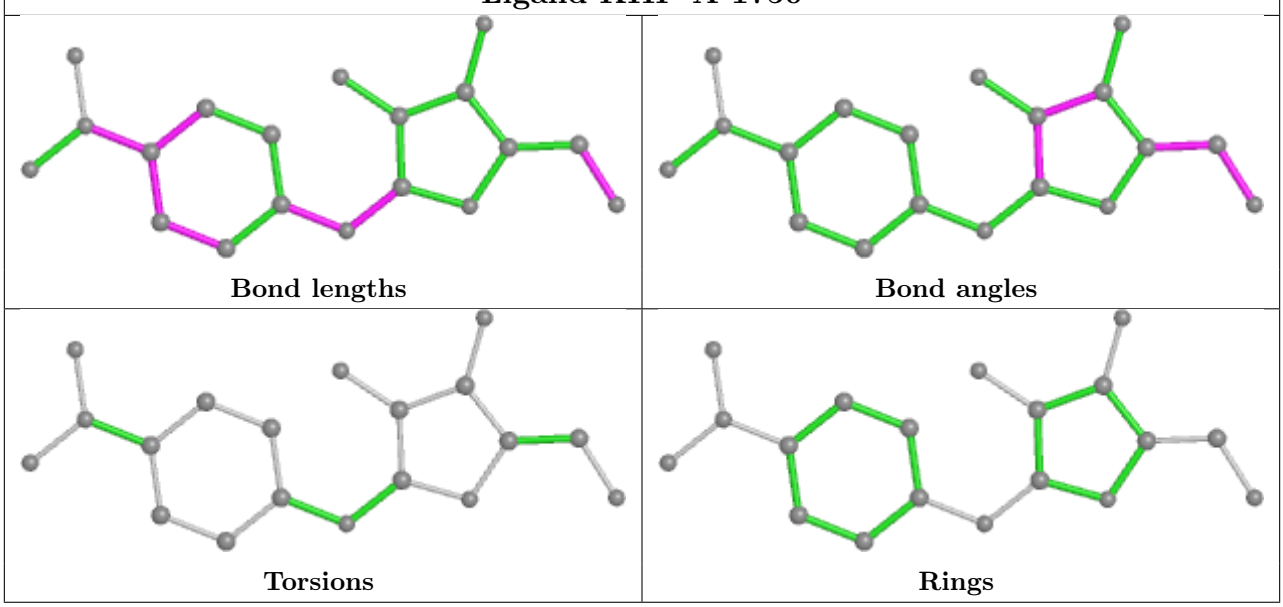
No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand KHP B 1751



Ligand KHP A 1750



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	497/502 (99%)	0.50	32 (6%)	19	18	8, 15, 30, 50	0
1	B	497/502 (99%)	0.18	22 (4%)	34	34	9, 15, 28, 38	0
All	All	994/1004 (99%)	0.34	54 (5%)	25	25	8, 15, 29, 50	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	443	PHE	6.4
1	A	482	ASP	6.3
1	A	441	ARG	6.0
1	A	444	GLU	5.4
1	A	14	PHE	4.9
1	A	252	ASN	4.7
1	A	445	ASP	4.7
1	A	178	GLY	4.7
1	A	417	GLU	4.5
1	B	441	ARG	4.3
1	B	180	TRP	4.2
1	B	482	ASP	4.0
1	A	180	TRP	4.0
1	B	14	PHE	3.9
1	A	391	VAL	3.9
1	A	483	ARG	3.9
1	A	457	ASP	3.8
1	B	49	GLU	3.6
1	B	480	LEU	3.5
1	A	251	ASP	3.5
1	B	466	GLN	3.4
1	A	466	GLN	3.3
1	A	434	LEU	3.3
1	A	402	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	170	TRP	3.0
1	A	442	SER	3.0
1	B	178	GLY	3.0
1	A	433	ALA	3.0
1	B	457	ASP	2.9
1	B	444	GLU	2.9
1	B	43	GLY	2.9
1	A	9	ILE	2.8
1	A	307	LEU	2.8
1	B	252	ASN	2.7
1	A	480	LEU	2.7
1	A	435	LEU	2.7
1	B	435	LEU	2.7
1	A	440	VAL	2.6
1	A	481	SER	2.5
1	A	49	GLU	2.5
1	A	5	LYS	2.5
1	B	481	SER	2.4
1	B	483	ARG	2.3
1	B	392	ILE	2.3
1	A	446	TYR	2.3
1	A	501	LYS	2.2
1	B	5	LYS	2.2
1	B	484	LYS	2.2
1	A	431	GLU	2.2
1	B	447	ARG	2.1
1	B	443	PHE	2.1
1	A	10	ILE	2.1
1	A	176	MET	2.1
1	B	445	ASP	2.1

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

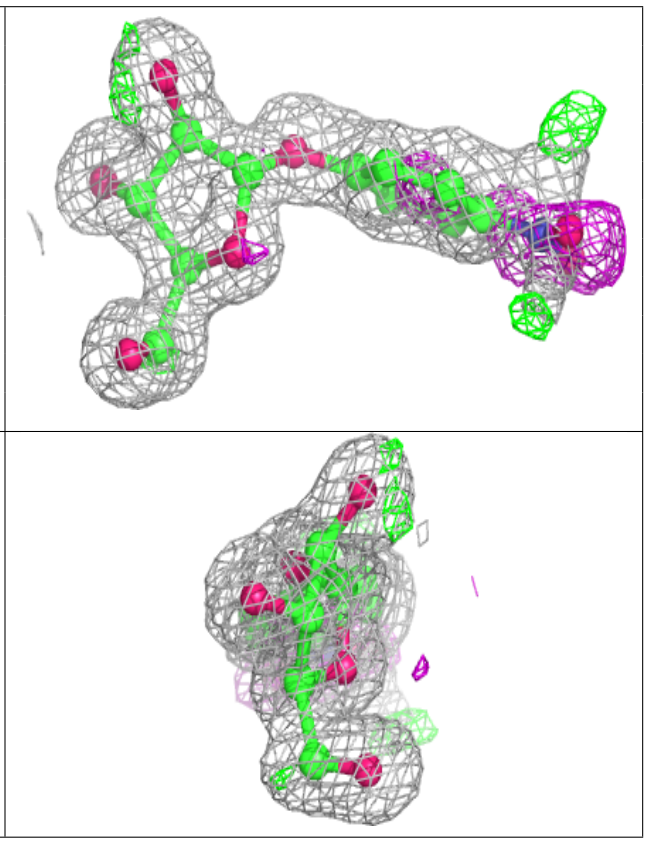
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	KHP	A	1750	19/19	0.91	0.14	10,13,42,43	0
2	KHP	B	1751	19/19	0.91	0.14	11,14,42,43	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

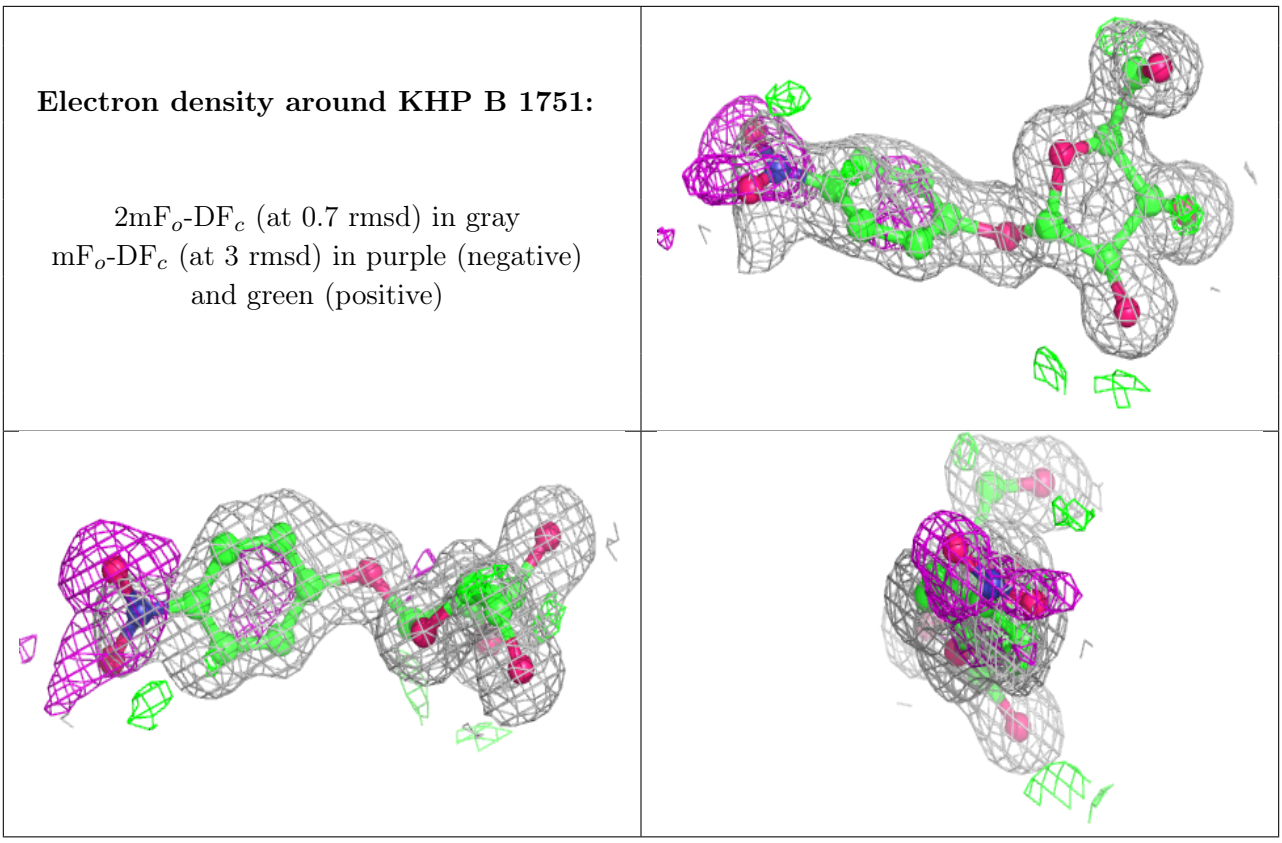
Electron density around KHP A 1750:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)



Electron density around KHP B 1751:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers (i)

There are no such residues in this entry.