Full wwPDB X-ray Structure Validation Report (i)

Aug 8, 2023 – 03:16 PM EDT

PDB ID	:	1P6B
Title	:	X-ray structure of phosphotriesterase, triple mutant H254G/H257W/L303T
Authors	:	Hill, C.M.; Li, W.; Thoden, J.B.; Holden, H.M.; Raushel, F.M.
Deposited on	:	2003-04-29
Resolution	:	1.90 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:

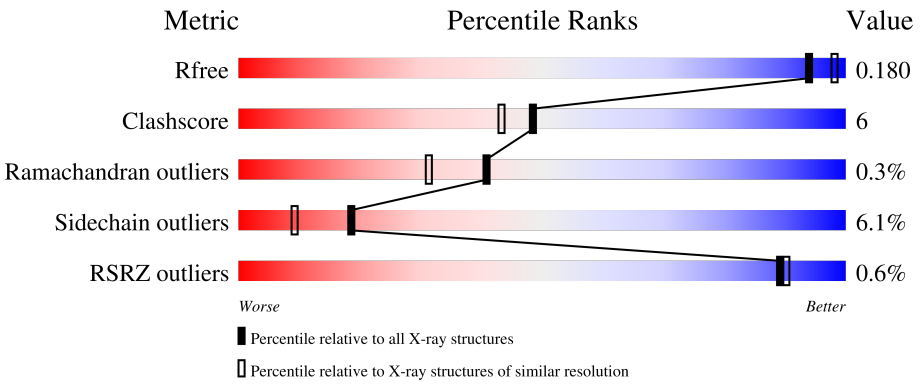
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	336	73% 21% 5% •
1	B	336	% 72% 23% • •

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5478 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 Parathion hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	330	2521	1595	447	472	7	0	0	0
1	B	331	2534	1603	448	475	8	0	3	0

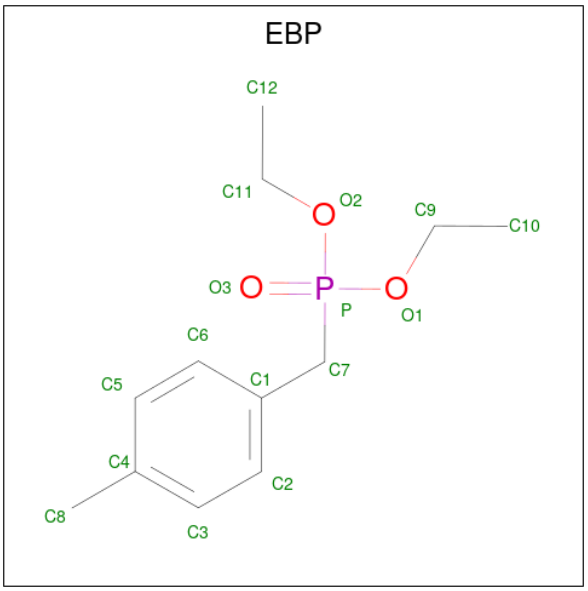
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	254	GLY	HIS	engineered mutation	UNP P0A433
A	257	TRP	HIS	engineered mutation	UNP P0A433
A	303	THR	LEU	engineered mutation	UNP P0A433
B	254	GLY	HIS	engineered mutation	UNP P0A433
B	257	TRP	HIS	engineered mutation	UNP P0A433
B	303	THR	LEU	engineered mutation	UNP P0A433

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

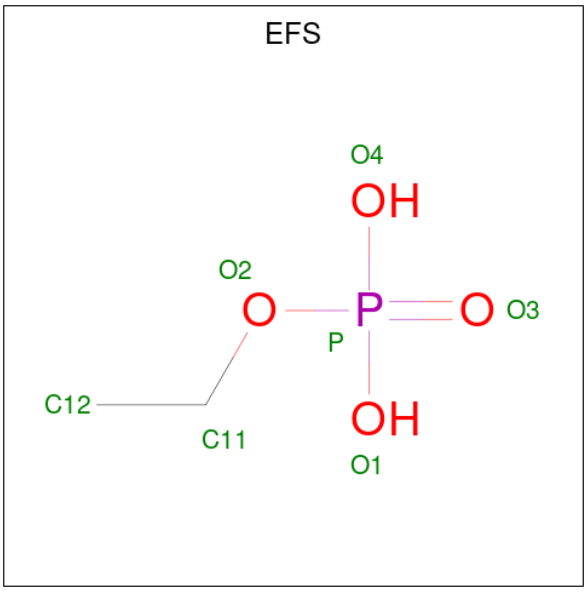
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
2	A	3	3	3	0	0
2	B	3	3	3	0	0

- Molecule 3 is DIETHYL 4-METHYLBENZYLPHOSPHONATE (three-letter code: EBP) (formula: $C_{12}H_{19}O_3P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
3	A	1	16	12	3	1	0	0
			Total	C	O	P		
3	B	1	16	12	3	1	0	0

- Molecule 4 is ETHYL DIHYDROGEN PHOSPHATE (three-letter code: EFS) (formula: $C_2H_7O_4P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
4	B	1	7	2	4	1	0	0

- Molecule 5 is water.

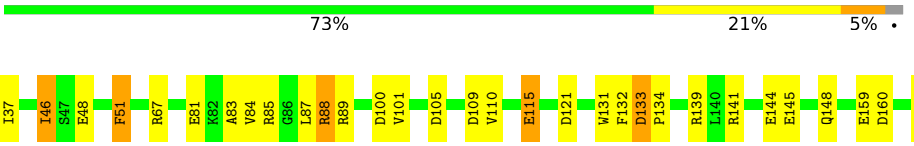
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	172	172	172	0	0
			Total	O		
5	B	206	206	206	0	0

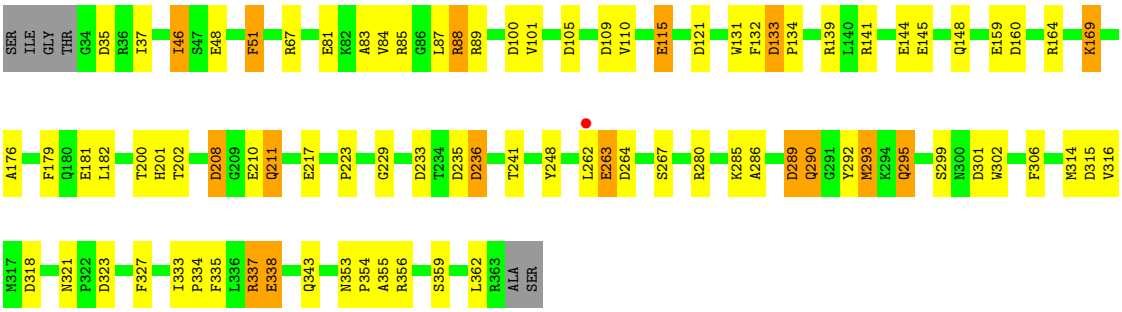
3 Residue-property plots (i)

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: Parathion hydrolase

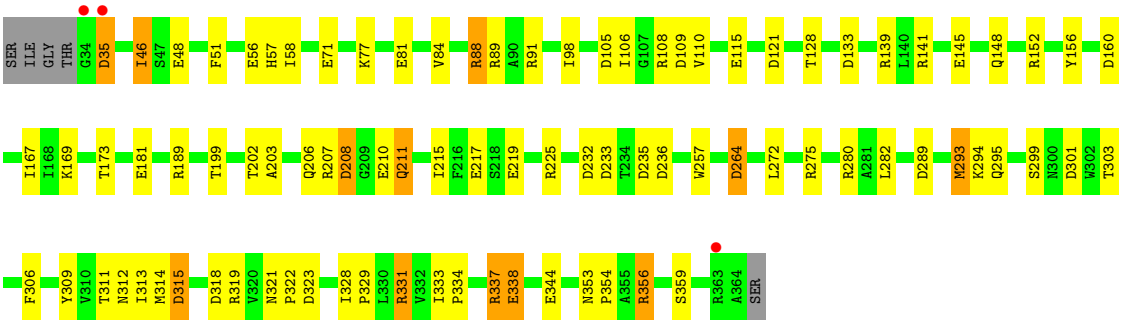
Chain A:





• Molecule 1: Parathion hydrolase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	130.40Å 91.90Å 69.70Å 90.00° 90.90° 90.00°	Depositor
Resolution (Å)	30.00 – 1.90 75.12 – 1.85	Depositor EDS
% Data completeness (in resolution range)	93.1 (30.00-1.90) 89.8 (75.12-1.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	2.08 (at 1.84Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.176 , 0.231 0.177 , 0.180	Depositor DCC
R_{free} test set	6306 reflections (9.96%)	wwPDB-VP
Wilson B-factor (Å^2)	17.5	Xtriage
Anisotropy	0.557	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 105.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.128 for -h,-k,l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	5478	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.28% 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: EBP, KCX, EFS, 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	0.87	11/2556 (0.4%)	1.43	45/3471 (1.3%)
1	B	0.85	12/2581 (0.5%)	1.38	49/3504 (1.4%)
All	All	0.86	23/5137 (0.4%)	1.41	94/6975 (1.3%)

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	338	GLU	CD-OE2	6.57	1.32	1.25
1	A	210	GLU	CD-OE2	6.08	1.32	1.25
1	A	217	GLU	CD-OE2	5.96	1.32	1.25
1	A	144	GLU	CD-OE2	5.83	1.32	1.25
1	B	219	GLU	CD-OE2	5.80	1.32	1.25
1	A	263	GLU	CD-OE2	5.68	1.31	1.25
1	A	181	GLU	CD-OE2	5.61	1.31	1.25
1	B	56	GLU	CD-OE2	5.61	1.31	1.25
1	B	210	GLU	CD-OE2	5.59	1.31	1.25
1	B	338	GLU	CD-OE2	5.59	1.31	1.25
1	B	344	GLU	CD-OE2	5.55	1.31	1.25
1	A	159	GLU	CD-OE2	5.48	1.31	1.25
1	B	71	GLU	CD-OE2	5.48	1.31	1.25
1	B	81	GLU	CD-OE2	5.47	1.31	1.25
1	A	145	GLU	CD-OE2	5.42	1.31	1.25
1	B	48	GLU	CD-OE2	5.41	1.31	1.25
1	B	181	GLU	CD-OE2	5.40	1.31	1.25
1	B	217	GLU	CD-OE2	5.40	1.31	1.25
1	B	115	GLU	CD-OE2	5.22	1.31	1.25
1	A	81	GLU	CD-OE2	5.21	1.31	1.25
1	B	145	GLU	CD-OE1	-5.14	1.20	1.25
1	A	48	GLU	CD-OE2	5.09	1.31	1.25
1	A	115	GLU	CD-OE2	5.09	1.31	1.25

All (94) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	337	ARG	NE-CZ-NH1	13.27	126.94	120.30
1	A	236	ASP	CB-CG-OD2	-11.28	108.15	118.30
1	A	85	ARG	NE-CZ-NH2	-11.06	114.77	120.30
1	A	85	ARG	NE-CZ-NH1	10.25	125.42	120.30
1	A	141	ARG	NE-CZ-NH1	10.05	125.32	120.30
1	B	139	ARG	NE-CZ-NH1	10.03	125.32	120.30
1	A	289	ASP	CB-CG-OD2	-9.64	109.62	118.30
1	B	105	ASP	CB-CG-OD2	-9.09	110.12	118.30
1	A	164	ARG	NE-CZ-NH1	8.70	124.65	120.30
1	B	225	ARG	NE-CZ-NH1	8.02	124.31	120.30
1	B	318	ASP	CB-CG-OD2	-8.01	111.09	118.30
1	A	35	ASP	CB-CG-OD1	7.87	125.38	118.30
1	B	233	ASP	CB-CG-OD1	7.72	125.25	118.30
1	A	356	ARG	NE-CZ-NH1	7.72	124.16	120.30
1	B	189	ARG	NE-CZ-NH1	7.72	124.16	120.30
1	A	236	ASP	CB-CG-OD1	7.60	125.14	118.30
1	A	318	ASP	CB-CG-OD2	-7.58	111.47	118.30
1	B	233	ASP	CB-CG-OD2	-7.55	111.51	118.30
1	B	35	ASP	CB-CG-OD2	-7.47	111.58	118.30
1	B	35	ASP	CB-CG-OD1	7.44	124.99	118.30
1	B	289	ASP	CB-CG-OD2	-7.43	111.61	118.30
1	A	35	ASP	CB-CG-OD2	-7.40	111.64	118.30
1	A	280	ARG	NE-CZ-NH1	7.26	123.93	120.30
1	A	160	ASP	CB-CG-OD2	-7.12	111.89	118.30
1	A	233	ASP	CB-CG-OD2	-7.12	111.90	118.30
1	B	301	ASP	CB-CG-OD1	7.10	124.69	118.30
1	A	89	ARG	NE-CZ-NH1	7.05	123.83	120.30
1	B	108	ARG	NE-CZ-NH1	6.97	123.78	120.30
1	B	91	ARG	NE-CZ-NH2	-6.96	116.82	120.30
1	A	289	ASP	CB-CG-OD1	6.94	124.55	118.30
1	B	225	ARG	NE-CZ-NH2	-6.92	116.84	120.30
1	B	121	ASP	CB-CG-OD2	-6.91	112.08	118.30
1	A	323	ASP	CB-CG-OD2	-6.90	112.09	118.30
1	B	208	ASP	CB-CG-OD2	-6.83	112.16	118.30
1	A	160	ASP	CB-CG-OD1	6.80	124.42	118.30
1	A	337	ARG	NE-CZ-NH2	-6.78	116.91	120.30
1	B	109	ASP	CB-CG-OD1	6.77	124.39	118.30
1	A	133	ASP	CB-CG-OD2	-6.77	112.21	118.30
1	A	67	ARG	NE-CZ-NH1	6.64	123.62	120.30
1	B	289	ASP	CB-CG-OD1	6.60	124.24	118.30
1	B	160	ASP	CB-CG-OD1	6.60	124.24	118.30
1	B	235	ASP	CB-CG-OD2	-6.56	112.39	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	232	ASP	CB-CG-OD2	-6.50	112.45	118.30
1	B	105	ASP	CB-CG-OD1	6.44	124.09	118.30
1	B	236	ASP	CB-CG-OD1	6.44	124.09	118.30
1	B	189	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	B	208	ASP	CB-CG-OD1	6.37	124.04	118.30
1	B	315	ASP	CB-CG-OD2	-6.37	112.56	118.30
1	B	141	ARG	NE-CZ-NH1	6.35	123.48	120.30
1	A	100	ASP	CB-CG-OD1	6.25	123.93	118.30
1	B	280	ARG	NE-CZ-NH1	6.25	123.43	120.30
1	B	232	ASP	CB-CG-OD1	6.25	123.93	118.30
1	A	139	ARG	NE-CZ-NH1	6.19	123.40	120.30
1	B	275	ARG	NE-CZ-NH1	6.15	123.37	120.30
1	A	109	ASP	CB-CG-OD1	6.12	123.81	118.30
1	B	152	ARG	NE-CZ-NH1	6.07	123.33	120.30
1	A	233	ASP	CB-CG-OD1	6.02	123.72	118.30
1	A	264	ASP	CB-CG-OD2	-6.00	112.90	118.30
1	A	356	ARG	NE-CZ-NH2	-5.99	117.31	120.30
1	A	315	ASP	CB-CG-OD2	-5.96	112.94	118.30
1	A	292	TYR	CA-CB-CG	-5.84	102.30	113.40
1	A	208	ASP	CB-CG-OD2	-5.80	113.08	118.30
1	A	89	ARG	NE-CZ-NH2	-5.80	117.40	120.30
1	A	301	ASP	CB-CG-OD1	5.79	123.51	118.30
1	B	264	ASP	CB-CG-OD2	-5.78	113.10	118.30
1	A	67	ARG	NE-CZ-NH2	-5.75	117.42	120.30
1	B	139	ARG	NE-CZ-NH2	-5.70	117.45	120.30
1	A	88	ARG	NE-CZ-NH1	5.69	123.14	120.30
1	B	235	ASP	CB-CG-OD1	5.61	123.35	118.30
1	A	235	ASP	CB-CG-OD2	-5.59	113.27	118.30
1	B	264	ASP	CB-CG-OD1	5.59	123.33	118.30
1	B	315	ASP	CB-CG-OD1	5.54	123.29	118.30
1	A	323	ASP	CB-CG-OD1	5.47	123.22	118.30
1	B	236	ASP	CB-CG-OD2	-5.44	113.41	118.30
1	A	335	PHE	N-CA-CB	-5.42	100.85	110.60
1	A	141	ARG	NE-CZ-NH2	-5.41	117.60	120.30
1	B	160	ASP	CB-CG-OD2	-5.40	113.44	118.30
1	A	105	ASP	CB-CG-OD2	-5.38	113.46	118.30
1	A	121	ASP	CB-CG-OD1	5.32	123.09	118.30
1	A	133	ASP	CB-CG-OD1	5.27	123.04	118.30
1	B	331	ARG	NE-CZ-NH1	5.27	122.93	120.30
1	B	89	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	B	318	ASP	CB-CG-OD1	5.26	123.03	118.30
1	B	91	ARG	NE-CZ-NH1	5.24	122.92	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	323	ASP	CB-CG-OD2	-5.21	113.61	118.30
1	B	128	THR	N-CA-CB	5.20	120.17	110.30
1	B	121	ASP	CB-CG-OD1	5.18	122.97	118.30
1	A	327	PHE	N-CA-CB	-5.18	101.27	110.60
1	B	203	ALA	N-CA-C	-5.14	97.12	111.00
1	B	356	ARG	NE-CZ-NH2	-5.13	117.74	120.30
1	A	264	ASP	CB-CG-OD1	5.07	122.86	118.30
1	B	133	ASP	CB-CG-OD2	-5.03	113.77	118.30
1	B	141	ARG	NE-CZ-NH2	-5.01	117.79	120.30
1	A	337	ARG	CD-NE-CZ	5.00	130.60	123.60

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	2521	0	2537	28	0
1	B	2534	0	2549	31	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	16	0	19	2	0
3	B	16	0	19	3	0
4	B	7	0	5	2	0
5	A	172	0	0	1	0
5	B	206	0	0	2	0
All	All	5478	0	5129	61	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 6.

All (61) 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:156:TYR:CG	3:B:8:EBP:H122	2.11	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:ILE:HB	1:A:334:PRO:HD3	1.64	0.78
1:A:293:MET:HG2	5:A:527:HOH:O	1.86	0.74
1:B:84:VAL:O	1:B:88:ARG:HG3	1.88	0.73
1:B:173:THR:HG22	4:B:9:EFS:O1	1.88	0.73
1:A:176:ALA:H	1:A:211:GLN:HE22	1.38	0.69
1:B:156:TYR:CD2	3:B:8:EBP:H122	2.28	0.69
1:A:46:ILE:HD13	1:A:359:SER:HB3	1.76	0.68
1:B:46:ILE:HD12	1:B:359:SER:OG	1.95	0.67
3:A:7:EBP:H6	3:A:7:EBP:H92	1.77	0.66
1:A:202:THR:HB	1:A:208:ASP:HB2	1.78	0.65
1:B:156:TYR:CD1	3:B:8:EBP:H122	2.32	0.65
1:B:293[B]:MET:HG2	5:B:483:HOH:O	1.98	0.63
1:B:294:LYS:O	1:B:356:ARG:NH2	2.32	0.62
1:A:285:LYS:NZ	1:A:289:ASP:OD2	2.33	0.62
1:B:337:ARG:HG2	5:B:425:HOH:O	2.00	0.61
1:A:176:ALA:H	1:A:211:GLN:NE2	1.98	0.61
1:A:46:ILE:CD1	1:A:359:SER:HB3	2.30	0.61
1:A:286:ALA:O	1:A:290:GLN:HG2	2.01	0.60
1:A:37:ILE:HD11	1:A:46:ILE:HD12	1.84	0.59
1:A:337:ARG:NH1	1:A:343:GLN:H	2.04	0.56
3:A:7:EBP:H6	3:A:7:EBP:C9	2.35	0.55
1:A:353:ASN:HB2	1:A:354:PRO:HD3	1.88	0.54
1:A:262:LEU:CD2	1:A:316:VAL:HG13	2.38	0.54
1:A:84:VAL:O	1:A:88:ARG:HG3	2.08	0.54
1:A:337:ARG:HH11	1:A:343:GLN:H	1.54	0.54
1:B:333:ILE:HB	1:B:334:PRO:HD3	1.90	0.54
1:B:353:ASN:HB2	1:B:354:PRO:HD3	1.91	0.53
1:A:302:TRP:CH2	1:A:321:ASN:HB3	2.44	0.52
1:B:328:ILE:HB	1:B:329:PRO:HD3	1.91	0.51
1:A:200:THR:O	1:A:229:GLY:HA3	2.11	0.51
1:A:169:KCX:OQ2	1:A:201:HIS:HB2	2.11	0.50
1:B:315:ASP:O	1:B:319:ARG:HD3	2.12	0.50
1:B:202:THR:HB	1:B:208:ASP:HB2	1.95	0.49
1:A:46:ILE:HD11	1:A:355:ALA:O	2.12	0.49
1:B:173:THR:HA	4:B:9:EFS:O1	2.14	0.48
1:B:309:TYR:HD2	1:B:313:ILE:HG12	1.80	0.46
1:B:206:GLN:O	1:B:207:ARG:HB2	2.15	0.46
1:B:282:LEU:HD23	1:B:282:LEU:HA	1.82	0.45
1:B:211:GLN:O	1:B:215:ILE:HD12	2.17	0.45
1:B:46:ILE:H	1:B:46:ILE:HG12	1.62	0.45
1:B:167:ILE:HD11	1:B:199:THR:HB	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:PHE:C	1:A:51:PHE:HD1	2.21	0.44
1:B:35:ASP:O	1:B:46:ILE:HG12	2.18	0.44
1:A:83:ALA:O	1:A:87:LEU:HG	2.18	0.44
1:B:57:HIS:O	1:B:303:THR:HA	2.18	0.44
1:A:51:PHE:C	1:A:51:PHE:CD1	2.90	0.44
1:B:106:ILE:HG22	1:B:106:ILE:O	2.18	0.43
1:B:311:THR:O	1:B:312:ASN:HB2	2.18	0.43
1:B:321:ASN:HA	1:B:322:PRO:HD2	1.81	0.43
1:A:179:PHE:O	1:A:182:LEU:HB2	2.20	0.42
1:A:295:GLN:HE21	1:A:295:GLN:HB3	1.59	0.42
1:B:257:TRP:CE3	1:B:272:LEU:HD23	2.55	0.42
1:B:338:GLU:O	1:B:338:GLU:HG2	2.19	0.42
1:A:262:LEU:HD21	1:A:316:VAL:HG13	2.02	0.41
1:A:131:TRP:CG	1:A:132:PHE:N	2.89	0.41
1:A:223:PRO:HA	1:A:248:TYR:CE1	2.56	0.40
1:A:133:ASP:N	1:A:134:PRO:CD	2.83	0.40
1:B:58:ILE:HD13	1:B:58:ILE:HA	1.73	0.40
1:B:98:ILE:O	1:B:98:ILE:HG13	2.21	0.40
1:B:328:ILE:HB	1:B:329:PRO:CD	2.52	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	327/336 (97%)	316 (97%)	9 (3%)	2 (1%)	25 15
1	B	331/336 (98%)	321 (97%)	10 (3%)	0	100 100
All	All	658/672 (98%)	637 (97%)	19 (3%)	2 (0%)	41 31

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	236	ASP
1	A	101	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	263/267 (98%)	246 (94%)	17 (6%)	17 8
1	B	266/267 (100%)	250 (94%)	16 (6%)	19 9
All	All	529/534 (99%)	496 (94%)	33 (6%)	18 9

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	ILE
1	A	51	PHE
1	A	110	VAL
1	A	115	GLU
1	A	148	GLN
1	A	211	GLN
1	A	241	THR
1	A	263	GLU
1	A	267	SER
1	A	290	GLN
1	A	293	MET
1	A	295	GLN
1	A	299	SER
1	A	306	PHE
1	A	314	MET
1	A	338	GLU
1	A	362	LEU
1	B	46	ILE
1	B	51	PHE
1	B	77	LYS
1	B	88	ARG
1	B	110	VAL

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Mol	Chain	Res	Type
1	B	148	GLN
1	B	211	GLN
1	B	264	ASP
1	B	293[A]	MET
1	B	293[B]	MET
1	B	295	GLN
1	B	299	SER
1	B	306	PHE
1	B	314	MET
1	B	331	ARG
1	B	337	ARG

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

Mol	Chain	Res	Type
1	A	211	GLN
1	A	212	GLN
1	A	295	GLN
1	B	212	GLN
1	B	312	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)

2 non-standard protein/DNA/RNA residues 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	$\#\|Z\| > 2$
1	KCX	A	169	2,1	9,11,12	3.09	2 (22%)	5,12,14	3.83	2 (40%)
1	KCX	B	169	2,1	9,11,12	2.70	2 (22%)	5,12,14	4.16	2 (40%)

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
1	KCX	A	169	2,1	-	0/9/10/12	-
1	KCX	B	169	2,1	-	1/9/10/12	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	169	KCX	CX-NZ	7.87	1.49	1.35
1	B	169	KCX	CX-NZ	5.88	1.45	1.35
1	B	169	KCX	OQ1-CX	5.43	1.31	1.21
1	A	169	KCX	OQ1-CX	4.47	1.30	1.21

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	169	KCX	OQ1-CX-NZ	-8.87	111.22	124.96
1	A	169	KCX	OQ1-CX-NZ	-8.05	112.47	124.96
1	A	169	KCX	CE-NZ-CX	-2.86	117.29	121.89
1	B	169	KCX	CE-NZ-CX	-2.47	117.92	121.89

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	169	KCX	CA-CB-CG-CD

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	169	KCX	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 9 ligands modelled in this entry, 6 are monoatomic - leaving 3 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	EBP	B	8	-	16,16,16	3.10	5 (31%)	21,21,21	1.36	1 (4%)
3	EBP	A	7	-	16,16,16	3.01	5 (31%)	21,21,21	1.35	1 (4%)
4	EFS	B	9	-	6,6,6	1.06	0	8,8,8	1.90	2 (25%)

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
3	EBP	B	8	-	-	1/13/13/13	0/1/1/1
3	EBP	A	7	-	-	1/13/13/13	0/1/1/1
4	EFS	B	9	-	-	2/4/4/4	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	8	EBP	P-C7	-10.98	1.62	1.79
3	A	7	EBP	P-C7	-10.43	1.63	1.79
3	B	8	EBP	C6-C5	2.89	1.44	1.38
3	A	7	EBP	C6-C5	2.62	1.43	1.38
3	A	7	EBP	C3-C2	2.51	1.43	1.38
3	A	7	EBP	C2-C1	2.47	1.44	1.38
3	B	8	EBP	C6-C1	2.45	1.44	1.38
3	B	8	EBP	C3-C2	2.17	1.42	1.38
3	A	7	EBP	C6-C1	2.12	1.43	1.38
3	B	8	EBP	C2-C1	2.10	1.43	1.38

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	7	EBP	O3-P-C7	-4.74	103.24	114.19
3	B	8	EBP	O3-P-C7	-4.46	103.88	114.19
4	B	9	EFS	O2-P-O3	3.13	115.27	106.47
4	B	9	EFS	O4-P-O2	-2.97	98.82	106.73

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	9	EFS	C11-O2-P-O4
4	B	9	EFS	C11-O2-P-O1
3	A	7	EBP	C9-O1-P-C7
3	B	8	EBP	C9-O1-P-C7

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	8	EBP	3	0
3	A	7	EBP	2	0
4	B	9	EFS	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)

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	329/336 (97%)	-0.62	1 (0%)	94	94	14, 23, 58, 97	0
1	B	330/336 (98%)	-0.65	3 (0%)	84	85	11, 22, 57, 100	0
All	All	659/672 (98%)	-0.63	4 (0%)	89	90	11, 23, 58, 100	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	262	LEU	3.8
1	B	363	ARG	2.4
1	B	34	GLY	2.1
1	B	35	ASP	2.1

6.2 Non-standard residues in protein, DNA, RNA chains (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
1	KCX	A	169	12/13	0.98	0.07	13,19,29,30	0
1	KCX	B	169	12/13	0.98	0.07	11,15,22,24	0

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	EBP	B	8	16/16	0.68	0.23	0,31,48,51	0
3	EBP	A	7	16/16	0.71	0.22	0,36,61,96	0
4	EFS	B	9	7/7	0.78	0.22	1,24,46,75	7
2	ZN	B	404	1/1	0.99	0.07	20,20,20,20	0
2	ZN	A	401	1/1	1.00	0.07	21,21,21,21	0
2	ZN	B	405	1/1	1.00	0.06	25,25,25,25	0
2	ZN	A	402	1/1	1.00	0.06	23,23,23,23	0
2	ZN	A	406	1/1	1.00	0.06	28,28,28,28	0
2	ZN	B	403	1/1	1.00	0.07	19,19,19,19	0

6.5 Other polymers (i)

There are no such residues in this entry.