

RETENTION TIMES OF INITIAL CALIBRATION

Lab Name:	<u>Alliance</u>	Contract:	<u>LANG01</u>
Lab Code:	<u>ACE</u>	SDG NO.:	<u>Q3310</u>
Instrument ID:	<u>ECD_L</u>	Calibration Date(s):	<u>10/07/2025</u> <u>10/07/2025</u>
		Calibration Times:	<u>10:33</u> <u>11:55</u>

GC Column: ZB-MR1 ID: 0.32 (mm)

LAB FILE ID:	RT 100 = <u>PL097526.D</u>	RT 075 = <u>PL097527.D</u>
RT 050 = <u>PL097528.D</u>	RT 025 = <u>PL097529.D</u>	RT 005 = <u>PL097530.D</u>

COMPOUND	RT 100	RT 075	RT 050	RT 025	RT 005	MEAN RT	RT WINDOW	
							FROM	TO
4,4'-DDD	6.66	6.66	6.66	6.66	6.66	6.66	6.56	6.76
4,4'-DDE	6.15	6.15	6.15	6.15	6.15	6.15	6.05	6.25
4,4'-DDT	6.98	6.98	6.98	6.98	6.98	6.97	6.87	7.07
Aldrin	5.23	5.23	5.23	5.23	5.23	5.23	5.13	5.33
alpha-BHC	3.97	3.97	3.97	3.97	3.97	3.97	3.87	4.07
alpha-Chlordane	5.98	5.98	5.98	5.98	5.98	5.98	5.88	6.08
beta-BHC	4.49	4.49	4.49	4.49	4.49	4.49	4.39	4.59
Decachlorobiphenyl	9.00	9.00	9.00	9.00	9.00	9.00	8.90	9.10
delta-BHC	4.73	4.73	4.73	4.73	4.73	4.73	4.63	4.83
Dieldrin	6.30	6.30	6.30	6.30	6.30	6.30	6.20	6.40
Endosulfan I	6.03	6.03	6.03	6.03	6.03	6.03	5.93	6.13
Endosulfan II	6.74	6.74	6.74	6.74	6.74	6.74	6.64	6.84
Endosulfan sulfate	7.10	7.10	7.10	7.10	7.10	7.10	7.00	7.20
Endrin	6.53	6.53	6.53	6.53	6.53	6.53	6.43	6.63
Endrin aldehyde	6.87	6.87	6.87	6.87	6.87	6.87	6.77	6.97
Endrin ketone	7.58	7.58	7.58	7.58	7.58	7.58	7.48	7.68
gamma-BHC (Lindane)	4.30	4.30	4.30	4.30	4.30	4.30	4.20	4.40
gamma-Chlordane	5.90	5.90	5.90	5.90	5.90	5.90	5.80	6.00
Heptachlor	4.89	4.89	4.89	4.89	4.89	4.89	4.79	4.99
Heptachlor epoxide	5.65	5.65	5.65	5.65	5.65	5.65	5.55	5.75
Methoxychlor	7.45	7.45	7.45	7.45	7.45	7.45	7.35	7.55
Tetrachloro-m-xylene	3.53	3.52	3.53	3.53	3.53	3.52	3.42	3.62



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GC Column: ZB-MR2 ID: 0.32 (mm)

LAB FILE ID:	RT 100 =	<u>PL097526.D</u>	RT 075 =	<u>PL097527.D</u>
RT 050 =	PL097528.D	RT 025 =	PL097529.D	RT 005 =
				PL097530.D

[illegible]