

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name:	<u>Alliance</u>	Contract:	<u>ALLI03</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3483</u>
Instrument ID:	<u>BNA_N</u>	Calibration Date/Time:	<u>12/01/2025 10:23</u>
Lab File ID:	<u>BN038222.D</u>	Init. Calib. Date(s):	<u>11/26/2025 11/27/2025</u>
EPA Sample No.:	<u>SSTDCCC0.4</u>	Init. Calib. Time(s):	<u>20:35 00:13</u>
GC Column:	<u>ZB-GR</u>	ID:	<u>0.25</u> (mm)

COMPOUND	RRF	RRF0.4	MIN RRF	%D	MAX%D
2-Methylnaphthalene-d10	0.568	0.586		3.2	20.0
Fluoranthene-d10	0.868	0.913		5.2	20.0
2-Fluorophenol	0.935	0.918		-1.8	20.0
Phenol-d6	1.168	1.147		-1.8	20.0
Nitrobenzene-d5	0.320	0.313		-2.2	20.0
Naphthalene	1.111	1.175		5.8	20.0
2-Methylnaphthalene	0.731	0.767		4.9	20.0
2-Fluorobiphenyl	1.545	1.642		6.3	20.0
Acenaphthylene	1.789	1.856		3.7	20.0
Acenaphthene	1.247	1.318		5.7	20.0
Fluorene	1.614	1.720		6.6	20.0
2,4,6-Tribromophenol	0.126	0.135		7.1	20.0
Phenanthrene	1.256	1.292		2.9	20.0
Anthracene	1.071	1.084		1.2	20.0
Fluoranthene	1.210	1.315		8.7	20.0
Pyrene	1.998	2.356		17.9	20.0
Terphenyl-d14	0.924	0.969		4.9	20.0
Benzo (a) anthracene	1.355	1.417		4.6	20.0
Chrysene	1.599	1.727		8.0	20.0
Benzo (b) fluoranthene	1.642	1.699		3.5	20.0
Benzo (k) fluoranthene	1.707	1.786		4.6	20.0
Benzo (a) pyrene	1.376	1.344		-2.3	20.0
Indeno (1,2,3-cd) pyrene	1.851	1.610		-13.0	20.0
Dibenzo (a,h) anthracene	1.401	1.136		-18.9	20.0
Benzo (g,h,i) perylene	1.674	1.675		0.1	20.0

All other compounds must meet a minimum RRF of 0.010.