

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM120216\
 Data File : BM008161.D
 Acq On : 02 Dec 2016 18:02
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.462

Quant Time: Dec 02 18:35:59 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 02 15:28:31 2016
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	95	-0.02
2 I	Naphthalene-d8	1.000	1.000	0.0	104	-0.01
3	Naphthalene	1.119	1.068	4.6	98	-0.01
4 SURR	2-Methylnaphthalene-d10	0.566	0.557	1.6	102	-0.01
5	2-Methylnaphthalene	0.721	0.702	2.6	101	0.00
6 I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
7	Acenaphthylene	1.810	1.766	2.4	99	0.00
8 C	Acenaphthene	1.377	1.307	5.1	98	0.00
9	Fluorene	1.583	1.426	9.9	92	0.00
10 I	Phenanthrene-d10	1.000	1.000	0.0	87	-0.02
11	Pentachlorophenol	0.140	0.104	25.7#	67	-0.02
12	Phenanthrene	1.246	1.154	7.4	81	0.00
13	Anthracene	1.170	1.201	-2.6	89	0.00
14 SURR	Fluoranthene-d10	1.132	1.023	9.6	79	0.00
15 C	Fluoranthene	1.608	1.277	20.6#	71	0.00
16 I	Chrysene-d12	1.000	1.000	0.0	60	0.00
17	Pyrene	1.389	1.650	-18.8	72	0.00
18	Benzo(a)anthracene	1.228	1.336	-8.8	68	0.00
19	Chrysene	1.527	1.349	11.7	55	0.00
20 I	Perylene-d12	1.000	1.000	0.0	63	0.00
21	Benzo(b)fluoranthene	1.596	1.545	3.2	62	0.00
22	Benzo(k)fluoranthene	1.647	1.451	11.9	57	0.00
23 C	Benzo(a)pyrene	1.495	1.475	1.3	63	0.00
24	Indeno(1,2,3-cd)pyrene	1.848	1.846	0.1	62	0.00
25	Dibenzo(a,h)anthracene	1.481	1.427	3.6	61	0.00
26	Benzo(g,h,i)perylene	1.640	1.620	1.2	61	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 1