

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM120216\
 Data File : BM008175.D
 Acq On : 03 Dec 2016 02:24
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.463

Quant Time: Dec 03 04:55:02 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 03 00:55:35 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	80	0.00
2 I	Naphthalene-d8	1.000	1.000	0.0	85	0.00
3	Naphthalene	1.119	1.079	3.6	82	0.00
4 SURR	2-Methylnaphthalene-d10	0.566	0.538	4.9	81	0.00
5	2-Methylnaphthalene	0.721	0.676	6.2	80	0.00
6 I	Acenaphthene-d10	1.000	1.000	0.0	77	0.00
7	Acenaphthylene	1.810	1.848	-2.1	79	0.00
8 C	Acenaphthene	1.377	1.363	1.0	78	0.00
9	Fluorene	1.583	1.499	5.3	74	0.00
10 I	Phenanthrene-d10	1.000	1.000	0.0	70	0.00
11	Pentachlorophenol	0.140	0.116	17.1	60	0.00
12	Phenanthrene	1.246	1.206	3.2	68	0.00
13	Anthracene	1.170	1.286	-9.9	77	0.00
14 SURR	Fluoranthene-d10	1.132	1.059	6.4	66	0.00
15 C	Fluoranthene	1.608	1.450	9.8	65	0.00
16 I	Chrysene-d12	1.000	1.000	0.0	58	0.00
17	Pyrene	1.389	1.573	-13.2	66	0.00
18	Benzo(a)anthracene	1.228	1.272	-3.6	62	0.00
19	Chrysene	1.527	1.410	7.7	55	0.00
20 I	Perylene-d12	1.000	1.000	0.0	62	0.00
21	Benzo(b)fluoranthene	1.596	1.422	10.9	56	0.00
22	Benzo(k)fluoranthene	1.647	1.565	5.0	61	0.00
23 C	Benzo(a)pyrene	1.495	1.439	3.7	61	0.00
24	Indeno(1,2,3-cd)pyrene	1.848	1.785	3.4	60	0.00
25	Dibenzo(a,h)anthracene	1.481	1.405	5.1	59	0.00
26	Benzo(g,h,i)perylene	1.640	1.594	2.8	60	0.00

(#) = Out of Range

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