

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121615\
 Data File : BM003594.D
 Acq On : 17 Dec 2015 02:55
 Operator : UM/IZ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.457

Quant Time: Dec 17 08:19:55 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM121515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 17 02:56:38 2015
 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	53	0.00
2 SURR	1,4-Dioxane-d8	0.171	0.179	-4.7	55	0.00
3	1,4-Dioxane	0.218	0.211	3.2	52	-0.01
4 I	Naphthalene-d8	1.000	1.000	0.0	60	0.00
5	Naphthalene	0.901	0.931	-3.3	61	0.00
6 SURR	2-Methylnaphthalene-d10	0.479	0.498	-4.0	61	0.00
7	2-Methylnaphthalene	0.605	0.670	-10.7	64	-0.01
8 I	Acenaphthene-d10	1.000	1.000	0.0	65	0.00
9	Acenaphthylene	1.528	1.499	1.9	64	0.00
10 C	Acenaphthene	1.232	1.328	-7.8	68	0.00
11	Fluorene	1.365	1.572	-15.2	73	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	70	0.00
13	Pentachlorophenol	0.057	0.036	36.8#	45	0.00
14	Phenanthrene	1.032	1.532	-48.4#	102	0.00
15	Anthracene	0.893	1.083	-21.3#	84	-0.02
16 SURR	Fluoranthene-d10	0.849	0.943	-11.1	75	0.00
17 C	Fluoranthene	1.120	1.659	-48.1#	100	0.00
18 I	Chrysene-d12	1.000	1.000	0.0	73	-0.01
19	Pyrene	1.080	1.436	-33.0#	92	-0.01
20	Benzo(a)anthracene	0.977	1.132	-15.9	83	-0.01
21	Chrysene	1.119	1.174	-4.9	80	-0.01
22 I	Perylene-d12	1.000	1.000	0.0	80	-0.01
23	Benzo(b)fluoranthene	1.365	1.492	-9.3	85	-0.01
24	Benzo(k)fluoranthene	1.237	1.153	6.8	76	-0.01
25 C	Benzo(a)pyrene	1.152	1.205	-4.6	83	-0.01
26	Indeno(1,2,3-cd)pyrene	1.335	1.259	5.7	80	-0.01
27	Dibenzo(a,h)anthracene	1.075	1.012	5.9	80	-0.01
28	Benzo(q,h,i)perylene	1.162	1.077	7.3	79	-0.01

(#) = Out of Range

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