

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121615\
 Data File : BM003579.D
 Acq On : 16 Dec 2015 16:06
 Operator : UM/IZ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.456

Quant Time: Dec 17 02:47:54 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 01:09:19 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	95	0.00
2 SURR	1,4-Dioxane-d8	0.171	0.171	0.0	94	0.00
3	1,4-Dioxane	0.218	0.207	5.0	91	0.00
4 I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
5	Naphthalene	0.901	0.917	-1.8	99	0.00
6 SURR	2-Methylnaphthalene-d10	0.479	0.487	-1.7	98	0.00
7	2-Methylnaphthalene	0.605	0.641	-6.0	101	0.00
8 I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
9	Acenaphthylene	1.528	1.548	-1.3	103	0.00
10 C	Acenaphthene	1.232	1.279	-3.8	103	0.00
11	Fluorene	1.365	1.441	-5.6	104	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
13	Pentachlorophenol	0.057	0.035	38.6#	63	0.00
14	Phenanthrene	1.032	1.077	-4.4	103	0.00
15	Anthracene	0.893	0.909	-1.8	101	0.00
16 SURR	Fluoranthene-d10	0.849	0.831	2.1	96	0.00
17 C	Fluoranthene	1.120	1.117	0.3	97	0.00
18 I	Chrysene-d12	1.000	1.000	0.0	84	0.00
19	Pyrene	1.080	1.301	-20.5#	96	0.00
20	Benzo(a)anthracene	0.977	1.008	-3.2	84	0.00
21	Chrysene	1.119	1.152	-2.9	89	0.00
22 I	Perylene-d12	1.000	1.000	0.0	89	0.00
23	Benzo(b)fluoranthene	1.365	1.313	3.8	82	0.00
24	Benzo(k)fluoranthene	1.237	1.336	-8.0	97	0.00
25 C	Benzo(a)pyrene	1.152	1.183	-2.7	91	0.00
26	Indeno(1,2,3-cd)pyrene	1.335	1.366	-2.3	96	0.00
27	Dibenzo(a,h)anthracene	1.075	1.102	-2.5	96	0.00
28	Benzo(q,h,i)perylene	1.162	1.206	-3.8	98	0.00

(#) = Out of Range

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