

Data Path : Z:\HPCHEM1\BNA M\Data\BM041116\
 Data File : BM004932.D
 Acq On : 11 Apr 2016 11:19
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.439

Quant Time: Apr 12 01:10:11 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM032516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 09 01:17:35 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	119	0.00
2 SURR	1,4-Dioxane-d8	0.180	0.194	-7.8	126	0.00
3	1,4-Dioxane	0.220	0.234	-6.4	124	0.00
4 I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
5	Naphthalene	0.966	0.967	-0.1	108	0.00
6 SURR	2-Methylnaphthalene-d10	0.558	0.500	10.4	96	0.00
7	2-Methylnaphthalene	0.692	0.664	4.0	101	0.00
8 I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
9	Acenaphthylene	1.840	1.636	11.1	85	0.00
10 C	Acenaphthene	1.302	1.356	-4.1	99	0.00
11	Fluorene	1.554	1.553	0.1	96	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
13	Pentachlorophenol	0.100	0.082	18.0	80	0.00
14	Phenanthrene	1.124	1.160	-3.2	100	0.00
15	Anthracene	1.062	0.977	8.0	89	0.00
16 SURR	Fluoranthene-d10	0.946	0.918	3.0	94	0.00
17 C	Fluoranthene	1.279	1.305	-2.0	99	0.00
18 I	Chrysene-d12	1.000	1.000	0.0	98	0.00
19	Pyrene	1.370	1.363	0.5	98	0.00
20	Benzo(a)anthracene	1.230	1.111	9.7	88	0.00
21	Chrysene	1.188	1.265	-6.5	102	0.00
22 I	Perylene-d12	1.000	1.000	0.0	88	0.00
23	Benzo(b)fluoranthene	1.348	1.525	-13.1	102	0.00
24	Benzo(k)fluoranthene	1.249	1.256	-0.6	86	0.00
25 C	Benzo(a)pyrene	1.232	1.271	-3.2	90	0.00
26	Indeno(1,2,3-cd)pyrene	1.140	1.237	-8.5	95	0.00
27	Dibenzo(a,h)anthracene	0.898	0.986	-9.8	97	0.00
28	Benzo(q,h,i)perylene	0.944	1.065	-12.8	100	0.00

(#) = Out of Range

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