

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032416\
 Data File : BM004769.D
 Acq On : 25 Mar 2016 04:19
 Operator : UM/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.435

Quant Time: Mar 25 04:33:37 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM032216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 25 03:18:48 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	84	0.00
2 SURR	1,4-Dioxane-d8	2.424	2.343	3.3	78	0.00
3	1,4-Dioxane	2.954	2.713	8.2	77	0.00
4 I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
5	Naphthalene	11.781	10.716	9.0	93	0.00
6 SURR	2-Methylnaphthalene-d10	5.695	5.705	-0.2	102	0.00
7	2-Methylnaphthalene	7.494	7.397	1.3	100	0.00
8 I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
9	Acenaphthylene	18.396	23.496	-27.7#	139	0.00
10 C	Acenaphthene	16.502	17.167	-4.0	111	0.00
11	Fluorene	17.439	19.085	-9.4	119	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	113	0.00
13	Pentachlorophenol	0.670	1.193	-78.1#	231#	0.00
14	Phenanthrene	13.789	13.974	-1.3	114	0.00
15	Anthracene	11.226	13.262	-18.1	137	0.00
16 SURR	Fluoranthene-d10	10.091	11.328	-12.3	129	0.00
17 C	Fluoranthene	14.171	16.243	-14.6	134	0.00
18 I	Chrysene-d12	1.000	1.000	0.0	6638#	-0.04
19	Pyrene	17.437	0.357	98.0#	132	-0.01
20	Benzo(a)anthracene	11.843	0.738	93.8#	398#	-0.04
21	Chrysene	16.264	0.243	98.5#	101	-0.01
22 I	Perylene-d12	1.000	1.000	0.0	8169#	-0.02
23	Benzo(b)fluoranthene	17.883	1.174	93.4#	536#	-0.01
24	Benzo(k)fluoranthene	18.118	0.535	97.0#	241#	-0.01
25 C	Benzo(a)pyrene	15.046	0.904	94.0#	490#	-0.02
26	Indeno(1,2,3-cd)pyrene	16.523	0.660	96.0#	322#	-0.01
27	Dibenzo(a,h)anthracene	12.860	0.324	97.5#	204#	-0.01
28	Benzo(q,h,i)perylene	15.728	0.608	96.1#	313#	-0.01

(#) = Out of Range

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