

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022816\
 Data File : BM004462.D
 Acq On : 28 Feb 2016 23:21
 Operator : SJ/UM
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Feb 29 06:41:25 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM022816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 29 05:57:18 2016
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	97	0.00
2	1,4-Dioxane	0.475	0.389	18.1	100	0.00
3	n-Nitrosodimethylamine	0.371	0.332	10.5	100	0.00
4 S	2-Fluorophenol	1.100	0.912	17.1	91	0.00
5 S	Phenol-d6	1.248	0.968	22.4	93	0.00
6	bis(2-Chloroethyl)ether	0.992	0.900	9.3	98	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
8 S	Nitrobenzene-d5	0.276	0.217	21.4	90	0.00
9	Nitrobenzene	0.307	0.250	18.6	92	0.00
10	Hexachlorobutadiene	0.204	0.176	13.7	95	0.00
11	2-Methylnaphthalene	0.761	0.697	8.4	102	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
13 S	2,4,6-Tribromophenol	0.198	0.118	40.4#	79	0.00
14 S	2-Fluorobiphenyl	1.525	1.266	17.0	99	0.00
15 I	Phenanthrene-d10	1.000	1.000	0.0	97	0.00
16	Hexachlorobenzene	0.278	0.251	9.7	101	0.00
17	Pentachlorophenol	0.103	0.061	40.8#	91	0.00
18	Fluoranthene	1.981	1.696	14.4	99	0.00
19 I	Chrysene-d12	1.000	1.000	0.0	100	0.00
20	Benzidine	0.257	0.335	-30.4#	158#	0.00
21	Pvrene	1.489	1.289	13.4	98	0.00
22 S	Terphenyl-d14	0.637	0.557	12.6	100	0.00
23	Benzo(a)anthracene	1.428	1.222	14.4	109	0.00
24	3,3'-Dichlorobenzidine	0.490	0.466	4.9	99	0.00
25	Chrysene	2.306	1.267	45.1#	100	0.00
26	Bis(2-ethylhexyl)phthalate	0.906	0.565	37.6#	75	0.00
27	Indeno(1,2,3-cd)pyrene	1.301	1.118	14.1	99	-0.01
28 I	Perylene-d12	1.000	1.000	0.0	101	-0.01
29	Benzo(b)fluoranthene	1.885	1.427	24.3	84	-0.01
30	Benzo(k)fluoranthene	1.625	1.453	10.6	111	0.00
31 C	Benzo(a)pyrene	1.466	1.291	11.9	100	0.00
32	Dibenzo(a,h)anthracene	1.309	1.091	16.7	99	0.00
33	Benzo(g,h,i)perylene	1.438	1.222	15.0	101	-0.01

(#) = Out of Range

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