

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122415\
 Data File : BM003774.D
 Acq On : 24 Dec 2015 12:39
 Operator : UM/SJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Dec 24 16:13:14 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM122415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 24 10:08:49 2015
 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	88	0.00
2	1,4-Dioxane	1.000	0.917	8.3	89	0.00
3	n-Nitrosodimethylamine	1.000	0.845	15.5	85	0.00
4 S	2-Fluorophenol	1.000	0.839	16.1	86	0.02
5 S	Phenol-d6	1.000	0.821	17.9	86	0.00
6	bis(2-Chloroethyl)ether	1.000	0.838	16.2	85	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	89	0.00
8 S	Nitrobenzene-d5	1.000	0.806	19.4	85	0.00
9	Nitrobenzene	1.000	0.800	20.0	84	0.00
10	Naphthalene	1.000	0.884	11.6	89	0.00
11	Hexachlorobutadiene	1.000	0.878	12.2	88	0.00
12	2-Methylnaphthalene	1.000	0.870	13.0	89	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	93	0.00
14 S	2,4,6-Tribromophenol	1.000	0.805	19.5	88	0.00
15 S	2-Fluorobiphenyl	1.000	0.844	15.6	90	0.00
16	Acenaphthylene	1.000	0.791	20.9	89	0.00
17	Acenaphthene	1.000	0.888	11.2	89	0.00
18	Fluorene	1.000	0.842	15.8	91	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	96	0.00
20	4-Bromophenyl-phenylether	1.000	1.197	-19.7	90	0.00
21	Hexachlorobenzene	1.000	1.224	-22.4	94	0.00
22	Pentachlorophenol	1.000	0.915	8.5	114	0.00
23	Phenanthrene	1.000	1.163	-16.3	96	0.00
24	Anthracene	1.000	0.900	10.0	93	0.00
25	Fluoranthene	1.000	0.972	2.8	98	0.00
26 I	Chrysene-d12	5.000	5.000	0.0	113	0.00
27	Benzydine	1.000	0.363	63.7#	27	0.00
28	Pvrene	1.000	0.815	18.5	101	0.00
29 S	Terphenyl-d14	1.000	0.842	15.8	103	0.00
30	Benzo(a)anthracene	1.000	0.806	19.4	110	0.00
31	3,3'-Dichlorobenzidine	1.000	0.827	17.3	92	0.00
32	Chrysene	1.000	0.920	8.0	108	0.00
33	Bis(2-ethylhexyl)phthalate	1.000	0.861	13.9	108	0.00
34	Indeno(1,2,3-cd)pyrene	1.000	1.035	-3.5	126	0.01
35 I	Perylene-d12	5.000	5.000	0.0	126	0.00
36	Benzo(b)fluoranthene	1.000	0.859	14.1	120	0.00
37	Benzo(k)fluoranthene	1.000	0.826	17.4	125	0.00
38 C	Benzo(a)pyrene	1.000	0.843	15.7	125	0.00
39	Dibenzo(a,h)anthracene	1.000	0.880	12.0	126	0.00
40	Benzo(a,h,i)perylene	1.000	0.876	12.4	124	0.00

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Compound	Amount	Calc.	%Dev Area	% Dev(min)

(#) = Out of Range				
SPCC's out = 0 CCC's out = 0				