

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE092315\
 Data File : BE090950.D
 Acq On : 23 Sep 2015 15:23
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Sep 24 03:58:09 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE091815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 18 18:17:34 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	113	0.00
2	1,4-Dioxane	0.795	0.885	-11.3	123	0.00
3	n-Nitrosodimethylamine	1.067	0.926	13.2	101	0.00
4 S	2-Fluorophenol	1.238	1.029	16.9	95	0.00
5 S	Phenol-d6	1.559	1.321	15.3	98	0.04
6	bis(2-Chloroethyl)ether	1.546	1.768	-14.4	141	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	115	0.00
8 S	Nitrobenzene-d5	0.345	0.275	20.3	93	0.01
9	Nitrobenzene	0.383	0.295	23.0	92	0.01
10	Naphthalene	1.126	1.106	1.8	114	0.00
11	Hexachlorobutadiene	0.166	0.185	-11.4	127	0.00
12	2-Methylnaphthalene	0.672	0.658	2.1	115	0.02
13 I	Acenaphthene-d10	1.000	1.000	0.0	124	0.00
14 S	2,4,6-Tribromophenol	0.196	0.147	25.0#	107	0.00
15 S	2-Fluorobiphenyl	1.418	1.381	2.6	120	0.00
16	Acenaphthylene	8.705	8.493	2.4	122	0.00
17	Acenaphthene	1.341	1.353	-0.9	127	0.00
18	Fluorene	1.705	1.653	3.0	124	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	120	0.00
20	4-Bromophenyl-phenylether	0.193	0.190	1.6	120	0.00
21	Hexachlorobenzene	0.906	0.976	-7.7	129	0.00
22	Pentachlorophenol	0.087	0.065	25.3#	99	0.04
23	Phenanthrene	1.292	1.208	6.5	115	0.00
24	Anthracene	1.144	1.170	-2.3	127	0.00
25	Fluoranthene	1.525	1.435	5.9	114	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	103	0.00
27	Pyrene	1.141	1.167	-2.3	113	0.00
28 S	Terphenyl-d14	0.638	0.633	0.8	112	0.01
29	Benzo(a)anthracene	1.235	1.058	14.3	92	0.00
30	Chrysene	1.352	1.558	-15.2	119	0.00
31	Bis(2-ethylhexyl)phthalate	0.515	0.442	14.2	100	0.00
32	Indeno(1,2,3-cd)pyrene	1.455	1.243	14.6	87	0.00
33 I	Perylene-d12	1.000	1.000	0.0	93	0.00
34	Benzo(b)fluoranthene	1.170	1.044	10.8	86	0.00
35	Benzo(k)fluoranthene	1.347	1.647	-22.3	115	0.00
36 C	Benzo(a)pyrene	1.113	1.184	-6.4	103	0.00
37	Dibenzo(a,h)anthracene	1.158	1.042	10.0	85	0.00
38	Benzo(g,h,i)perylene	1.240	1.163	6.2	89	0.00

(#) = Out of Range

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