

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070115\
 Data File : BE090042.D
 Acq On : 1 Jul 2015 12:45
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Jul 02 06:34:21 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 30 07:01:01 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	100	-0.02
2	1,4-Dioxane	0.686	0.589	14.1	89	-0.01
3	n-Nitrosodimethylamine	0.930	0.760	18.3	85	-0.02
4 S	2-Fluorophenol	1.149	1.040	9.5	91	-0.01
5 S	Phenol-d6	1.608	1.473	8.4	93	-0.01
6 I	Naphthalene-d8	1.000	1.000	0.0	100	-0.01
7 S	Nitrobenzene-d5	0.397	0.339	14.6	87	-0.01
8	Nitrobenzene	0.421	0.361	14.3	88	-0.01
9	Naphthalene	1.130	1.041	7.9	93	-0.01
10	Hexachlorobutadiene	0.178	0.169	5.1	96	-0.01
11	2-Methylnaphthalene	0.755	0.675	10.6	91	-0.01
12 I	Acenaphthene-d10	1.000	1.000	0.0	94	-0.01
13 S	2,4,6-Tribromophenol	0.147	0.114	22.4	75	0.00
14 S	2-Fluorobiphenyl	1.503	1.360	9.5	87	-0.01
15	Acenaphthylene	9.019	8.318	7.8	90	0.00
16	Acenaphthene	1.300	1.227	5.6	90	-0.01
17	Fluorene	1.768	1.638	7.4	90	-0.02
18 I	Phenanthrene-d10	1.000	1.000	0.0	89	-0.02
19	4-Bromophenyl-phenylether	0.213	0.190	10.8	81	-0.02
20	Hexachlorobenzene	0.885	0.810	8.5	84	-0.02
21	Pentachlorophenol	0.086	0.054	37.2#	59	0.00
22	Phenanthrene	1.281	1.137	11.2	81	-0.02
23	Anthracene	1.172	1.076	8.2	83	0.00
24	Fluoranthene	1.431	1.289	9.9	82	-0.01
25 I	Chrysene-d12	1.000	1.000	0.0	87	-0.01
26	Pyrene	1.192	1.106	7.2	82	-0.01
27 S	Terphenyl-d14	0.833	0.741	11.0	77	-0.01
28	Benzo(a)anthracene	1.249	1.140	8.7	81	-0.01
29	Chrysene	1.297	1.163	10.3	79	-0.01
30	Bis(2-ethylhexyl)phthalate	0.398	0.353	11.3	84	-0.01
31	Indeno(1,2,3-cd)pyrene	1.243	1.182	4.9	84	-0.03
32 I	Perylene-d12	1.000	1.000	0.0	84	-0.02
33	Benzo(b)fluoranthene	1.103	1.134	-2.8	87	-0.01
34	Benzo(k)fluoranthene	1.232	1.219	1.1	83	-0.01
35 C	Benzo(a)pyrene	1.019	1.031	-1.2	85	-0.02
36	Dibenzo(a,h)anthracene	1.046	1.016	2.9	83	-0.03
37	Benzo(g,h,i)perylene	1.060	1.027	3.1	82	-0.03

(#) = Out of Range

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