

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081415\
 Data File : BE090530.D
 Acq On : 14 Aug 2015 21:00
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Aug 17 04:19:26 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE081415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 14 16:52:42 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	105	0.00
2	1,4-Dioxane	0.658	0.677	-2.9	102	0.00
3	n-Nitrosodimethylamine	0.863	0.870	-0.8	107	0.00
4 S	2-Fluorophenol	1.446	1.686	-16.6	121	0.00
5 S	Phenol-d6	2.125	2.421	-13.9	124	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
7 S	Nitrobenzene-d5	0.398	0.439	-10.3	119	0.00
8	Nitrobenzene	0.413	0.457	-10.7	121	0.00
9	Naphthalene	1.136	1.141	-0.4	104	0.00
10	Hexachlorobutadiene	0.167	0.164	1.8	101	0.00
11	2-Methylnaphthalene	0.740	0.760	-2.7	108	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
13 S	2,4,6-Tribromophenol	0.192	0.236	-22.9	141	0.00
14 S	2-Fluorobiphenyl	1.493	1.500	-0.5	105	0.00
15	Acenaphthylene	9.132	9.370	-2.6	109	0.00
16	Acenaphthene	1.374	1.389	-1.1	107	0.00
17	Fluorene	1.684	1.681	0.2	105	0.02
18 I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
19	4-Bromophenyl-phenylether	0.204	0.224	-9.8	111	0.00
20	Hexachlorobenzene	0.906	0.929	-2.5	102	0.00
21	Pentachlorophenol	0.079	0.116	-46.8#	192#	0.00
22	Phenanthrene	1.312	1.306	0.5	100	0.00
23	Anthracene	1.208	1.295	-7.2	110	0.00
24	Fluoranthene	1.396	1.493	-6.9	109	0.00
25 I	Chrysene-d12	1.000	1.000	0.0	101	0.00
26	Pyrene	1.430	1.515	-5.9	109	0.00
27 S	Terphenyl-d14	0.914	0.988	-8.1	110	0.00
28	Benzo(a)anthracene	1.316	1.353	-2.8	109	0.00
29	Chrysene	1.344	1.365	-1.6	103	0.00
30	Bis(2-ethylhexyl)phthalate	0.725	1.066	-47.0#	180#	0.00
31	Indeno(1,2,3-cd)pyrene	1.266	1.412	-11.5	121	0.00
32 I	Perylene-d12	1.000	1.000	0.0	108	0.00
33	Benzo(b)fluoranthene	1.184	1.285	-8.5	120	0.00
34	Benzo(k)fluoranthene	1.342	1.360	-1.3	108	0.00
35 C	Benzo(a)pyrene	1.108	1.188	-7.2	120	0.00
36	Dibenzo(a,h)anthracene	1.069	1.172	-9.6	125	0.00
37	Benzo(g,h,i)perylene	1.169	1.266	-8.3	119	0.00

(#) = Out of Range

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