

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081715\
 Data File : BE090546.D
 Acq On : 17 Aug 2015 23:09
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Aug 18 04:36:55 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	115	0.00
2	1,4-Dioxane	0.658	0.647	1.7	107	0.00
3	n-Nitrosodimethylamine	0.863	0.815	5.6	110	0.00
4 S	2-Fluorophenol	1.446	1.537	-6.3	121	0.00
5 S	Phenol-d6	2.125	2.193	-3.2	124	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	118	0.00
7 S	Nitrobenzene-d5	0.398	0.396	0.5	123	0.00
8	Nitrobenzene	0.413	0.409	1.0	125	0.00
9	Naphthalene	1.136	1.115	1.8	117	0.00
10	Hexachlorobutadiene	0.167	0.164	1.8	116	0.00
11	2-Methylnaphthalene	0.740	0.752	-1.6	123	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	121	0.00
13 S	2,4,6-Tribromophenol	0.192	0.203	-5.7	141	0.00
14 S	2-Fluorobiphenyl	1.493	1.506	-0.9	123	0.00
15	Acenaphthylene	9.132	9.201	-0.8	125	0.00
16	Acenaphthene	1.374	1.367	0.5	122	0.00
17	Fluorene	1.684	1.670	0.8	121	0.02
18 I	Phenanthrene-d10	1.000	1.000	0.0	119	0.00
19	4-Bromophenyl-phenylether	0.204	0.208	-2.0	124	0.00
20	Hexachlorobenzene	0.906	0.907	-0.1	119	0.00
21	Pentachlorophenol	0.079	0.075	5.1	150#	0.00
22	Phenanthrene	1.312	1.268	3.4	117	0.00
23	Anthracene	1.208	1.201	0.6	122	0.00
24	Fluoranthene	1.396	1.363	2.4	119	0.00
25 I	Chrysene-d12	1.000	1.000	0.0	116	0.00
26	Pyrene	1.430	1.446	-1.1	120	0.00
27 S	Terphenyl-d14	0.914	0.944	-3.3	122	0.00
28	Benzo(a)anthracene	1.316	1.290	2.0	120	0.00
29	Chrysene	1.344	1.342	0.1	117	0.00
30	Bis(2-ethylhexyl)phthalate	0.725	0.821	-13.2	160#	0.00
31	Indeno(1,2,3-cd)pyrene	1.266	1.279	-1.0	127	0.00
32 I	Perylene-d12	1.000	1.000	0.0	122	0.00
33	Benzo(b)fluoranthene	1.184	1.200	-1.4	126	0.00
34	Benzo(k)fluoranthene	1.342	1.366	-1.8	123	0.00
35 C	Benzo(a)pyrene	1.108	1.124	-1.4	127	0.00
36	Dibenzo(a,h)anthracene	1.069	1.073	-0.4	129	0.00
37	Benzo(g,h,i)perylene	1.169	1.196	-2.3	126	0.00

(#) = Out of Range

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