

Data Path : Z:\HPCHEM1\BNA E\DATA\BE080515\
 Data File : BE090376.D
 Acq On : 6 Aug 2015 1:18
 Operator : TP/UM
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Aug 06 13:07:16 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 06 12:28:08 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	116	0.00
2	1,4-Dioxane	0.687	0.795	-15.7	125	0.00
3	n-Nitrosodimethylamine	0.773	0.787	-1.8	119	0.00
4 S	2-Fluorophenol	0.750	0.925	-23.3	154#	0.00
5 S	Phenol-d6	1.241	1.354	-9.1	137	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	117	0.00
7 S	Nitrobenzene-d5	0.331	0.336	-1.5	125	0.00
8	Nitrobenzene	0.349	0.343	1.7	128	0.00
9	Naphthalene	1.104	1.109	-0.5	118	0.00
10	Hexachlorobutadiene	0.191	0.189	1.0	115	0.00
11	2-Methylnaphthalene	0.640	0.683	-6.7	130	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	118	0.00
13 S	2,4,6-Tribromophenol	0.095	0.104	-9.5	166#	0.00
14 S	2-Fluorobiphenyl	1.375	1.463	-6.4	124	0.00
15	Acenaphthylene	8.098	8.491	-4.9	125	0.00
16	Acenaphthene	1.259	1.256	0.2	118	0.00
17	Fluorene	1.635	1.643	-0.5	117	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	112	0.00
19	4-Bromophenyl-phenylether	0.194	0.210	-8.2	123	0.00
20	Hexachlorobenzene	1.005	1.019	-1.4	113	0.00
21	Pentachlorophenol	0.044	0.047	-6.8	192#	0.00
22	Phenanthrene	1.251	1.245	0.5	112	0.00
23	Anthracene	1.028	1.093	-6.3	122	-0.02
24	Fluoranthene	1.231	1.352	-9.8	127	0.00
25 I	Chrysene-d12	1.000	1.000	0.0	117	0.00
26	Pyrene	1.084	1.169	-7.8	131	0.00
27 S	Terphenyl-d14	0.737	0.792	-7.5	127	0.00
28	Benzo(a)anthracene	1.108	1.106	0.2	130	0.00
29	Chrysene	1.339	1.353	-1.0	115	0.00
30	Bis(2-ethylhexyl)phthalate	0.311	0.296	4.8	155#	0.00
31	Indeno(1,2,3-cd)pyrene	0.942	0.945	-0.3	142	0.00
32 I	Perylene-d12	1.000	1.000	0.0	123	0.00
33	Benzo(b)fluoranthene	0.918	1.043	-13.6	148	0.00
34	Benzo(k)fluoranthene	1.392	1.331	4.4	114	0.00
35 C	Benzo(a)pyrene	0.850	0.919	-8.1	141	0.00
36	Dibenzo(a,h)anthracene	0.766	0.793	-3.5	147	0.00
37	Benzo(g,h,i)perylene	0.923	0.982	-6.4	135	0.00

(#) = Out of Range

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