

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081415\
 Data File : BE090523.D
 Acq On : 14 Aug 2015 16:43
 Operator : UM/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE081415

Quant Time: Aug 17 04:17:49 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	136	0.00
2	1,4-Dioxane	0.658	0.641	2.6	125	0.00
3	n-Nitrosodimethylamine	0.863	0.830	3.8	133	0.00
4 S	2-Fluorophenol	1.446	1.370	5.3	128	0.00
5 S	Phenol-d6	2.125	2.044	3.8	137	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	135	0.00
7 S	Nitrobenzene-d5	0.398	0.387	2.8	138	0.00
8	Nitrobenzene	0.413	0.399	3.4	140	0.00
9	Naphthalene	1.136	1.096	3.5	131	0.00
10	Hexachlorobutadiene	0.167	0.165	1.2	134	0.00
11	2-Methylnaphthalene	0.740	0.715	3.4	134	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	137	0.00
13 S	2,4,6-Tribromophenol	0.192	0.176	8.3	139	-0.02
14 S	2-Fluorobiphenyl	1.493	1.456	2.5	134	0.00
15	Acenaphthylene	9.132	8.864	2.9	136	0.00
16	Acenaphthene	1.374	1.339	2.5	135	0.00
17	Fluorene	1.684	1.669	0.9	137	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	137	0.00
19	4-Bromophenyl-phenylether	0.204	0.199	2.5	137	0.00
20	Hexachlorobenzene	0.906	0.877	3.2	133	-0.02
21	Pentachlorophenol	0.079	0.067	15.2	154#	0.00
22	Phenanthrene	1.312	1.230	6.3	131	0.00
23	Anthracene	1.208	1.166	3.5	137	0.00
24	Fluoranthene	1.396	1.330	4.7	134	0.00
25 I	Chrysene-d12	1.000	1.000	0.0	138	0.00
26	Pyrene	1.430	1.384	3.2	136	0.00
27 S	Terphenyl-d14	0.914	0.892	2.4	136	0.00
28	Benzo(a)anthracene	1.316	1.244	5.5	137	0.00
29	Chrysene	1.344	1.302	3.1	134	0.00
30	Bis(2-ethylhexyl)phthalate	0.725	0.617	14.9	142	0.00
31	Indeno(1,2,3-cd)pyrene	1.266	1.245	1.7	146	0.00
32 I	Perylene-d12	1.000	1.000	0.0	139	0.00
33	Benzo(b)fluoranthene	1.184	1.132	4.4	136	0.00
34	Benzo(k)fluoranthene	1.342	1.367	-1.9	141	0.00
35 C	Benzo(a)pyrene	1.108	1.082	2.3	140	0.00
36	Dibenzo(a,h)anthracene	1.069	1.066	0.3	146	0.00
37	Benzo(g,h,i)perylene	1.169	1.154	1.3	140	0.00

(#) = Out of Range

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