

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
 Data File : BE089990.D
 Acq On : 29 Jun 2015 19:08
 Operator : TP/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE062915

Quant Time: Jun 30 07:02:35 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	130	0.00
2	1,4-Dioxane	0.686	0.577	15.9	114	0.00
3	n-Nitrosodimethylamine	0.930	0.780	16.1	114	0.00
4 S	2-Fluorophenol	1.149	0.961	16.4	109	0.00
5 S	Phenol-d6	1.608	1.385	13.9	114	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	134	0.00
7 S	Nitrobenzene-d5	0.397	0.334	15.9	114	0.00
8	Nitrobenzene	0.421	0.355	15.7	116	0.00
9	Naphthalene	1.130	1.029	8.9	124	0.00
10	Hexachlorobutadiene	0.178	0.165	7.3	126	0.00
11	2-Methylnaphthalene	0.755	0.677	10.3	122	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	135	0.00
13 S	2,4,6-Tribromophenol	0.147	0.114	22.4	107	0.00
14 S	2-Fluorobiphenyl	1.503	1.327	11.7	123	0.00
15	Acenaphthylene	9.019	8.136	9.8	126	0.00
16	Acenaphthene	1.300	1.224	5.8	129	0.00
17	Fluorene	1.768	1.630	7.8	128	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	134	0.00
19	4-Bromophenyl-phenylether	0.213	0.189	11.3	122	0.00
20	Hexachlorobenzene	0.885	0.821	7.2	130	0.00
21	Pentachlorophenol	0.086	0.058	32.6#	96	0.00
22	Phenanthrene	1.281	1.180	7.9	127	0.00
23	Anthracene	1.172	1.082	7.7	126	0.00
24	Fluoranthene	1.431	1.320	7.8	126	0.00
25 I	Chrysene-d12	1.000	1.000	0.0	137	0.00
26	Pyrene	1.192	1.090	8.6	128	0.00
27 S	Terphenyl-d14	0.833	0.737	11.5	121	0.00
28	Benzo(a)anthracene	1.249	1.145	8.3	128	0.00
29	Chrysene	1.297	1.160	10.6	124	0.00
30	Bis(2-ethylhexyl)phthalate	0.398	0.245	38.4#	92	0.00
31	Indeno(1,2,3-cd)pyrene	1.243	1.134	8.8	127	0.00
32 I	Perylene-d12	1.000	1.000	0.0	137	0.00
33	Benzo(b)fluoranthene	1.103	1.046	5.2	131	0.00
34	Benzo(k)fluoranthene	1.232	1.124	8.8	126	0.00
35 C	Benzo(a)pyrene	1.019	0.925	9.2	125	0.00
36	Dibenzo(a,h)anthracene	1.046	0.938	10.3	125	0.00
37	Benzo(g,h,i)perylene	1.060	0.952	10.2	125	0.00

(#) = Out of Range

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