

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030516\
 Data File : BM004566.D
 Acq On : 05 Mar 2016 21:08
 Operator : SJ/UM
 Sample : SSTDICV001
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM030516

Quant Time: Mar 07 07:58:25 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM030516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 07 07:54:56 2016
 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	119	0.00
2	1,4-Dioxane	0.592	0.493	16.7	120	0.00
3	n-Nitrosodimethylamine	0.385	0.349	9.4	120	0.00
4 S	2-Fluorophenol	1.184	1.062	10.3	119	0.00
5 S	Phenol-d6	1.421	1.303	8.3	121	0.00
6	bis(2-Chloroethyl)ether	1.203	1.151	4.3	123	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	117	0.00
8 S	Nitrobenzene-d5	0.267	0.247	7.5	121	0.00
9	Nitrobenzene	0.298	0.270	9.4	121	0.00
10	Naphthalene	1.237	1.149	7.1	132	0.00
11	Hexachlorobutadiene	0.216	0.203	6.0	122	0.00
12	2-Methylnaphthalene	0.797	0.746	6.4	123	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	116	0.00
14 S	2,4,6-Tribromophenol	0.189	0.162	14.3	120	0.00
15 S	2-Fluorobiphenyl	1.639	1.534	6.4	123	0.00
16	Acenaphthylene	2.566	2.343	8.7	121	0.00
17	Acenaphthene	1.547	1.410	8.9	125	0.00
18	Fluorene	2.075	1.914	7.8	122	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	120	0.00
20	Hexachlorobenzene	0.260	0.208	20.0	121	0.00
21	Pentachlorophenol	0.027	0.015	44.4#	113	0.00
22	Phenanthrene	1.149	1.034	10.0	127	0.00
23	Anthracene	1.134	0.964	15.0	126	0.00
24	Fluoranthene	1.460	1.291	11.6	123	0.00
25 I	Chrysene-d12	1.000	1.000	0.0	117	0.00
26	Benzidine	0.234	0.311	-32.9#	213#	-0.02
27	Pyrene	1.868	1.699	9.0	125	0.00
28 S	Terphenyl-d14	0.788	0.706	10.4	125	0.00
29	Benzo(a)anthracene	1.620	1.487	8.2	123	0.00
30	3,3'-Dichlorobenzidine	0.500	0.514	-2.8	126	0.00
31	Chrysene	1.689	1.552	8.1	126	-0.02
32	Bis(2-ethylhexyl)phthalate	1.215	0.952	21.6	113	0.00
33	Indeno(1,2,3-cd)pyrene	1.555	1.405	9.6	124	0.00
34 I	Perylene-d12	1.000	1.000	0.0	122	0.00
35 C	Benzo(a)pyrene	1.534	1.440	6.1	126	-0.01
36	Dibenzo(a,h)anthracene	1.278	1.149	10.1	125	0.00
37	Benzo(g,h,i)perylene	1.412	1.270	10.1	125	0.00

(#) = Out of Range

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