

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020216\
 Data File : BE091384.D
 Acq On : 2 Feb 2016 19:09
 Operator : UM/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE020216

Quant Time: Feb 03 00:35:50 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE020216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 03 00:27:53 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	101	0.00
2 S	2-Fluorophenol	0.896	0.973	-8.6	125	0.00
3 S	Phenol-d6	1.352	1.537	-13.7	121	0.00
4 I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
5 S	Nitrobenzene-d5	0.241	0.263	-9.1	131	0.00
6	Naphthalene	1.128	1.083	4.0	113	0.00
7	2-Methylnaphthalene	0.609	0.614	-0.8	119	0.00
8 I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
9 S	2,4,6-Tribromophenol	0.153	0.152	0.7	134	0.00
10 S	2-Fluorobiphenyl	1.301	1.235	5.1	114	0.00
11	Acenaphthylene	7.133	6.895	3.3	121	0.00
12	Acenaphthene	1.373	1.266	7.8	114	0.00
13	Fluorene	1.772	1.710	3.5	115	0.00
14 I	Phenanthrene-d10	1.000	1.000	0.0	107	0.00
15	Phenanthrene	1.313	1.230	6.3	114	0.00
16	Anthracene	1.100	1.079	1.9	117	0.00
17	Fluoranthene	1.203	1.360	-13.1	140	0.00
18 I	Chrysene-d12	1.000	1.000	0.0	112	0.00
19	Pyrene	1.083	1.115	-3.0	130	0.00
20 S	Terphenyl-d14	0.740	0.768	-3.8	128	0.00
21	Benzo(a)anthracene	0.926	1.016	-9.7	146	0.00
22	Chrysene	1.371	1.270	7.4	119	0.00
23	Indeno(1,2,3-cd)pyrene	0.929	1.072	-15.4	141	0.00
24 I	Perylene-d12	1.000	1.000	0.0	116	0.00
25	Benzo(b)fluoranthene	1.048	1.119	-6.8	135	0.00
26	Benzo(k)fluoranthene	1.187	1.194	-0.6	131	0.00
27 C	Benzo(a)pyrene	0.939	1.050	-11.8	144	0.00
28	Dibenzo(a,h)anthracene	0.827	0.996	-20.4	154#	0.00
29	Benzo(g,h,i)perylene	1.055	1.110	-5.2	136	0.00

(#) = Out of Range

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