

Data Path : V:\HPCHEM1\BNA E\DATA\BE070916\
 Data File : BE091996.D
 Acq On : 9 Jul 2016 13:27
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 11 04:51:47 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE070516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 05 16:14:58 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	97	0.00
2	1,4-Dioxane	0.678	0.662	2.4	91	0.00
3	n-Nitrosodimethylamine	0.768	0.692	9.9	93	-0.02
4 S	2-Fluorophenol	1.196	1.045	12.6	89	0.00
5 S	Phenol-d6	1.662	1.396	16.0	90	0.00
6	bis(2-Chloroethyl)ether	1.449	1.367	5.7	98	-0.02
7	n-Nitroso-di-n-propylamine	1.387	1.153	16.9	86	-0.02
8 I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
9 S	Nitrobenzene-d5	0.341	0.311	8.8	92	-0.02
10	Nitrobenzene	0.370	0.325	12.2	93	0.00
11	bis(2-Chloroethoxy)methane	0.418	0.428	-2.4	88	0.00
12	Naphthalene	1.164	1.130	2.9	94	0.00
13	Hexachlorobutadiene	0.177	0.165	6.8	86	0.00
14	2-Methylnaphthalene	0.748	0.703	6.0	91	-0.01
15 I	Acenaphthene-d10	1.000	1.000	0.0	90	0.00
16 S	2,4,6-Tribromophenol	0.143	0.130	9.1	96	0.00
17 S	2-Fluorobiphenyl	1.220	1.297	-6.3	93	0.00
18	Acenaphthylene	6.276	6.350	-1.2	94	0.00
19	2,6-Dinitrotoluene	1.093	0.979	10.4	112	0.00
20	Acenaphthene	1.367	1.299	5.0	84	0.00
21	2,4-Dinitrotoluene	1.115	1.335	-19.7	165#	0.00
22	Fluorene	1.442	1.469	-1.9	90	0.00
23	4-Chlorophenyl-phenylether	0.719	0.736	-2.4	88	0.00
24 I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
25	4-Bromophenyl-phenylether	0.217	0.193	11.1	88	0.00
26	Hexachlorobenzene	0.216	0.203	6.0	94	0.00
27	Pentachlorophenol	0.078	0.047	39.7#	105	0.00
28	Phenanthrene	1.303	1.252	3.9	93	0.00
29	Anthracene	1.218	1.108	9.0	93	0.00
30	Fluoranthene	5.599	5.209	7.0	94	0.00
31 I	Chrysene-d12	1.000	1.000	0.0	120	0.00
32	Benzidine	0.333	0.148	55.6#	63	0.00
33	Pyrene	3.640	2.970	18.4	97	0.00
34 S	Terphenyl-d14	0.590	0.527	10.7	106	0.00
35	Benzo(a)anthracene	0.837	0.836	0.1	121	0.00
36	3,3'-Dichlorobenzidine	0.257	0.151	41.2#	70	0.00
37	Chrysene	1.187	1.271	-7.1	129	0.00
38	Bis(2-ethylhexyl)phthalate	0.970	0.616	36.5#	83	-0.03
39	Indeno(1,2,3-cd)pyrene	4.657	3.832	17.7	96	-0.02
40 I	Perylene-d12	1.000	1.000	0.0	110	0.00
41	Benzo(b)fluoranthene	1.381	1.367	1.0	110	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.335	1.155	13.5	96	0.00
43 C	Benzo(a)pyrene	1.304	1.166	10.6	99	0.00
44	Dibenzo(a,h)anthracene	1.227	1.045	14.8	96	0.00
45	Benzo(a,h,i)perylene	5.004	4.296	14.1	95	0.00

(#) = Out of Range

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