

Data Path : Z:\HPCHEM1\BNA E\DATA\BE070716\
 Data File : BE091984.D
 Acq On : 7 Jul 2016 17:09
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 08 01:13:06 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	99	0.00
2	1,4-Dioxane	0.678	0.666	1.8	94	0.00
3	n-Nitrosodimethylamine	0.768	0.685	10.8	95	0.00
4 S	2-Fluorophenol	1.196	0.974	18.6	86	0.00
5 S	Phenol-d6	1.662	1.304	21.5	86	0.00
6	bis(2-Chloroethyl)ether	1.449	1.328	8.4	98	0.00
7	n-Nitroso-di-n-propylamine	1.387	1.137	18.0	87	-0.02
8 I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
9 S	Nitrobenzene-d5	0.341	0.291	14.7	87	-0.02
10	Nitrobenzene	0.370	0.304	17.8	88	0.00
11	bis(2-Chloroethoxy)methane	0.418	0.430	-2.9	90	0.00
12	Naphthalene	1.164	1.113	4.4	93	0.00
13	Hexachlorobutadiene	0.177	0.169	4.5	90	0.00
14	2-Methylnaphthalene	0.748	0.693	7.4	91	-0.01
15 I	Acenaphthene-d10	1.000	1.000	0.0	93	0.00
16 S	2,4,6-Tribromophenol	0.143	0.102	28.7#	78	0.00
17 S	2-Fluorobiphenyl	1.220	1.223	-0.2	90	0.00
18	Acenaphthylene	6.276	5.673	9.6	86	0.00
19	2,6-Dinitrotoluene	1.093	0.810	25.9#	95	0.00
20	Acenaphthene	1.367	1.318	3.6	87	0.00
21	2,4-Dinitrotoluene	1.115	0.653	41.4#	83	0.00
22	Fluorene	1.442	1.362	5.5	85	0.00
23	4-Chlorophenyl-phenylether	0.719	0.699	2.8	86	0.00
24 I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
25	4-Bromophenyl-phenylether	0.217	0.198	8.8	85	0.00
26	Hexachlorobenzene	0.216	0.207	4.2	90	0.00
27	Pentachlorophenol	0.078	0.041	47.4#	87	0.00
28	Phenanthrene	1.303	1.273	2.3	90	0.00
29	Anthracene	1.218	1.097	9.9	87	0.00
30	Fluoranthene	5.599	5.581	0.3	96	0.00
31 I	Chrysene-d12	1.000	1.000	0.0	97	0.00
32	Benzidine	0.333	0.172	48.3#	59	0.02
33	Pyrene	3.640	3.570	1.9	94	0.00
34 S	Terphenyl-d14	0.590	0.603	-2.2	98	0.00
35	Benzo(a)anthracene	0.837	0.739	11.7	86	0.00
36	3,3'-Dichlorobenzidine	0.257	0.188	26.8#	70	0.00
37	Chrysene	1.187	1.190	-0.3	98	0.00
38	Bis(2-ethylhexyl)phthalate	0.970	0.693	28.6#	75	0.00
39	Indeno(1,2,3-cd)pyrene	4.657	4.295	7.8	87	0.00
40 I	Perylene-d12	1.000	1.000	0.0	99	0.00
41	Benzo(b)fluoranthene	1.381	1.229	11.0	89	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.335	1.240	7.1	93	0.00
43 C	Benzo(a)pyrene	1.304	1.135	13.0	87	0.00
44	Dibenzo(a,h)anthracene	1.227	1.046	14.8	86	0.00
45	Benzo(a,h,i)perylene	5.004	4.361	12.8	87	0.00

(#) = Out of Range

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