

Data Path : Z:\HPCHEM1\BNA_E\Data\BE070717\
 Data File : BE093448.D
 Acq On : 7 Jul 2017 18:12
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 08 05:35:15 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE070217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 03 06:40:23 2017
 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	126	0.00
2	1,4-Dioxane	0.847	0.725	14.4	108	0.00
3	n-Nitrosodimethylamine	0.444	0.422	5.0	151#	0.00
4 S	2-Fluorophenol	1.207	1.156	4.2	124	0.00
5 S	Phenol-d6	1.780	1.802	-1.2	136	0.00
6	bis(2-Chloroethyl)ether	1.339	1.309	2.2	126	-0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	140	0.00
8 S	Nitrobenzene-d5	0.256	0.285	-11.3	169#	0.00
9	Naphthalene	4.190	4.067	2.9	140	0.00
10	Hexachlorobutadiene	0.159	0.148	6.9	133	-0.02
11 SURR	2-Methylnaphthalene-d10	0.638	0.653	-2.4	148	0.00
12	2-Methylnaphthalene	0.704	0.714	-1.4	148	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	141	0.00
14 S	2,4,6-Tribromophenol	0.126	0.085	32.5#	105	0.00
15 S	2-Fluorobiphenyl	1.553	1.539	0.9	145	0.00
16	Acenaphthylene	1.666	1.617	2.9	145	0.00
17	Acenaphthene	1.220	1.195	2.0	144	0.00
18	Fluorene	1.675	1.514	9.6	132	-0.02
19 I	Phenanthrene-d10	1.000	1.000	0.0	87	0.00
20	4-Bromophenyl-phenylether	0.219	0.332	-51.6#	142	0.00
21	Hexachlorobenzene	0.237	0.309	-30.4#	126	0.00
22	Pentachlorophenol	0.035	0.054	-54.3#	164#	0.00
23	Phenanthrene	1.454	1.583	-8.9	103	0.00
24	Anthracene	0.891	0.888	0.3	91	0.00
25 SURR	Fluoranthene-d10	4.823	5.488	-13.8	108	0.00
26	Fluoranthene	1.367	1.554	-13.7	108	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	101	0.00
28	Pyrene	1.124	1.158	-3.0	106	0.00
29 S	Terphenyl-d14	0.714	0.721	-1.0	104	0.00
30	Benzo(a)anthracene	1.080	1.031	4.5	101	0.00
31	Chrysene	1.130	1.116	1.2	101	0.00
32	Bis(2-ethylhexyl)phthalate	1.555	1.511	2.8	120	-0.01
33	Indeno(1,2,3-cd)pyrene	0.992	1.033	-4.1	108	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	106	0.00
35	Benzo(b)fluoranthene	1.292	1.215	6.0	102	0.00
36	Benzo(k)fluoranthene	1.278	1.237	3.2	108	0.00
37 C	Benzo(a)pyrene	1.132	1.092	3.5	108	0.00
38	Dibenzo(a,h)anthracene	1.099	1.072	2.5	107	-0.01
39	Benzo(g,h,i)perylene	1.110	1.060	4.5	106	0.00

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(#) = Out of Range	SPCC's out = 0 CCC's out = 0			