

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013117\
 Data File : BE092689.D
 Acq On : 31 Jan 2017 16:27
 Operator : SJ/MA
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Feb 01 10:34:49 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE011617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 17 10:14:01 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	156#	0.00
2	1,4-Dioxane	0.775	0.704	9.2	134	0.00
3	n-Nitrosodimethylamine	0.902	0.567	37.1#	112	0.02
4 S	2-Fluorophenol	1.116	1.004	10.0	165#	0.00
5 S	Phenol-d6	1.368	1.339	2.1	185#	-0.02
6	bis(2-Chloroethyl)ether	1.453	1.288	11.4	164#	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	162#	0.00
8 S	Nitrobenzene-d5	0.325	0.283	12.9	170#	-0.02
9	Naphthalene	1.044	1.012	3.1	170#	-0.02
10	Hexachlorobutadiene	0.154	0.160	-3.9	177#	-0.02
11	2-Methylnaphthalene	0.648	0.617	4.8	172#	-0.01
12 I	Acenaphthene-d10	1.000	1.000	0.0	159#	0.00
13 S	2,4,6-Tribromophenol	0.126	0.123	2.4	195#	0.02
14 S	2-Fluorobiphenyl	1.230	1.218	1.0	170#	-0.01
15	Acenaphthylene	6.890	6.912	-0.3	180#	-0.02
16	Acenaphthene	1.105	1.112	-0.6	169#	0.00
17	Fluorene	1.387	1.345	3.0	172#	-0.01
18 I	Phenanthrene-d10	1.000	1.000	0.0	166#	0.00
19	Hexachlorobenzene	0.189	0.170	10.1	143	0.00
20	Pentachlorophenol	0.051	0.038	25.5#	135	0.28
21	Phenanthrene	1.104	1.035	6.3	161#	0.00
22	Anthracene	1.018	0.991	2.7	181#	-0.02
23	Fluoranthene	4.895	4.993	-2.0	187#	-0.02
24 I	Chrysene-d12	1.000	1.000	0.0	151#	0.00
25	Pyrene	4.929	5.192	-5.3	191#	-0.02
26 S	Terphenyl-d14	0.707	0.721	-2.0	179#	0.00
27	Benzo(a)anthracene	4.926	5.237	-6.3	186#	0.00
28	Chrysene	5.542	6.037	-8.9	177#	0.00
29	Bis(2-ethylhexyl)phthalate	0.913	1.107	-21.2	256#	0.00
30	Indeno(1,2,3-cd)pyrene	5.987	6.573	-9.8	189#	-0.02
31 I	Perylene-d12	1.000	1.000	0.0	162#	0.00
32	Benzo(b)fluoranthene	1.210	1.209	0.1	176#	-0.01
33	Benzo(k)fluoranthene	1.215	1.278	-5.2	198#	-0.01
34 C	Benzo(a)pyrene	1.136	1.185	-4.3	195#	-0.01
35	Dibenzo(a,h)anthracene	1.112	1.151	-3.5	189#	-0.03
36	Benzo(g,h,i)perylene	4.699	4.841	-3.0	184#	-0.02

(#) = Out of Range

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