

Data Path : Z:\HPCHEM1\BNA_E\Data\BE072017\
 Data File : BE093496.D
 Acq On : 20 Jul 2017 11:27
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 20 18:15:07 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	108	0.00
2	1,4-Dioxane	0.847	0.722	14.8	92	0.00
3	n-Nitrosodimethylamine	0.444	0.455	-2.5	139	0.00
4 S	2-Fluorophenol	1.207	1.314	-8.9	121	0.00
5 S	Phenol-d6	1.780	2.195	-23.3	142	0.01
6	bis(2-Chloroethyl)ether	1.339	1.339	0.0	111	-0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	133	0.00
8 S	Nitrobenzene-d5	0.256	0.300	-17.2	169#	0.00
9	Naphthalene	4.190	3.942	5.9	129	0.00
10	Hexachlorobutadiene	0.159	0.148	6.9	126	-0.02
11 SURR	2-Methylnaphthalene-d10	0.638	0.638	0.0	138	0.00
12	2-Methylnaphthalene	0.704	0.699	0.7	138	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	138	0.00
14 S	2,4,6-Tribromophenol	0.126	0.152	-20.6	183#	0.00
15 S	2-Fluorobiphenyl	1.553	1.471	5.3	135	0.00
16	Acenaphthylene	1.666	1.630	2.2	142	0.00
17	Acenaphthene	1.220	1.129	7.5	133	0.00
18	Fluorene	1.675	1.443	13.9	123	-0.02
19 I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
20	4-Bromophenyl-phenylether	0.219	0.252	-15.1	131	0.00
21	Hexachlorobenzene	0.237	0.245	-3.4	121	0.00
22	Pentachlorophenol	0.035	0.047	-34.3#	170#	0.00
23	Phenanthrene	1.454	1.319	9.3	104	0.00
24	Anthracene	0.891	0.863	3.1	108	0.00
25 SURR	Fluoranthene-d10	4.823	4.415	8.5	105	0.00
26	Fluoranthene	1.367	1.223	10.5	103	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	98	0.00
28	Pyrene	1.124	1.155	-2.8	103	0.00
29 S	Terphenyl-d14	0.714	0.712	0.3	99	0.00
30	Benzo(a)anthracene	1.080	1.097	-1.6	105	0.00
31	Chrysene	1.130	1.055	6.6	93	0.00
32	Bis(2-ethylhexyl)phthalate	1.555	2.796	-79.8#	215#	-0.01
33	Indeno(1,2,3-cd)pyrene	0.992	1.018	-2.6	103	0.00
34 I	Perylene-d12	1.000	1.000	0.0	107	0.00
35	Benzo(b)fluoranthene	1.292	1.169	9.5	99	0.00
36	Benzo(k)fluoranthene	1.278	1.095	14.3	96	0.00
37 C	Benzo(a)pyrene	1.132	1.056	6.7	105	0.00
38	Dibenzo(a,h)anthracene	1.099	1.011	8.0	102	0.00
39	Benzo(g,h,i)perylene	1.110	1.005	9.5	101	0.00

Data Path : Z:\HPCHEM1\BNA_E\Data\BE072017\
Data File : BE093496.D
Acq On : 20 Jul 2017 11:27
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Jul 20 18:15:07 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)

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