

Data Path : Z:\HPCHEM1\BNA_E\Data\BE080517\
 Data File : BE093691.D
 Acq On : 6 Aug 2017 1:05
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 06 03:46:32 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 02 10:58:31 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	123	0.00
2	1,4-Dioxane	0.648	0.596	8.0	113	0.00
3	n-Nitrosodimethylamine	0.607	0.635	-4.6	136	0.00
4 S	2-Fluorophenol	1.192	1.334	-11.9	143	0.00
5 S	Phenol-d6	1.794	2.054	-14.5	145	0.00
6	bis(2-Chloroethyl)ether	1.508	1.523	-1.0	127	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	129	0.00
8 S	Nitrobenzene-d5	0.305	0.344	-12.8	156#	0.00
9	Naphthalene	4.630	4.558	1.6	128	0.00
10	Hexachlorobutadiene	0.172	0.174	-1.2	134	-0.02
11 SURR	2-Methylnaphthalene-d10	0.632	0.645	-2.1	132	0.00
12	2-Methylnaphthalene	0.766	0.766	0.0	130	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	129	0.00
14 S	2,4,6-Tribromophenol	0.093	0.107	-15.1	181#	0.00
15 S	2-Fluorobiphenyl	1.601	1.568	2.1	125	-0.02
16	Acenaphthylene	1.912	2.037	-6.5	142	0.00
17	Acenaphthene	1.349	1.323	1.9	128	0.00
18	Fluorene	1.803	1.721	4.5	122	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	123	0.00
20	4-Bromophenyl-phenylether	0.126	0.125	0.8	110	0.00
21	Hexachlorobenzene	0.129	0.141	-9.3	124	0.00
22	Pentachlorophenol	0.055	0.108	-96.4#	253#	0.00
23	Phenanthrene	0.880	0.880	0.0	120	0.00
24	Anthracene	1.203	1.334	-10.9	129	0.00
25 SURR	Fluoranthene-d10	4.541	4.945	-8.9	123	0.00
26	Fluoranthene	1.401	1.480	-5.6	120	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	119	-0.01
28	Pyrene	1.292	1.323	-2.4	123	0.00
29 S	Terphenyl-d14	0.719	0.714	0.7	117	0.00
30	Benzo(a)anthracene	1.176	1.307	-11.1	137	0.00
31	Chrysene	1.311	1.288	1.8	117	0.00
32	Bis(2-ethylhexyl)phthalate	1.603	3.578	-123.2#	310#	-0.01
33	Indeno(1,2,3-cd)pyrene	1.187	1.275	-7.4	127	0.00
34 I	Perylene-d12	1.000	1.000	0.0	128	0.00
35	Benzo(b)fluoranthene	1.427	1.389	2.7	128	0.00
36	Benzo(k)fluoranthene	1.459	1.342	8.0	122	0.00
37 C	Benzo(a)pyrene	1.293	1.289	0.3	132	0.00
38	Dibenzo(a,h)anthracene	1.248	1.200	3.8	125	0.00
39	Benzo(g,h,i)perylene	1.266	1.208	4.6	125	0.00

Data Path : Z:\HPCHEM1\BNA_E\Data\BE080517\
Data File : BE093691.D
Acq On : 6 Aug 2017 1:05
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Aug 06 03:46:32 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 02 10:58:31 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			