

Data Path : Z:\HPCHEM1\BNA E\Data\BE021318\
 Data File : BE095670.D
 Acq On : 12 Feb 2018 20:19
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Feb 13 04:23:04 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE020618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 20:27:34 2018
 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	97	0.00
2	1,4-Dioxane	0.662	0.614	7.3	90	0.00
3	n-Nitrosodimethylamine	0.833	0.823	1.2	95	0.00
4 S	2-Fluorophenol	1.207	1.285	-6.5	102	0.00
5 S	Phenol-d6	1.671	1.894	-13.3	109	-0.01
6	bis(2-Chloroethyl)ether	1.538	1.515	1.5	93	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
8 S	Nitrobenzene-d5	0.424	0.433	-2.1	101	0.00
9	Naphthalene	4.646	4.643	0.1	97	-0.02
10	Hexachlorobutadiene	0.249	0.274	-10.0	99	0.00
11 SURR	2-Methylnaphthalene-d10	0.662	0.679	-2.6	100	0.00
12	2-Methylnaphthalene	0.789	0.809	-2.5	99	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
14 S	2,4,6-Tribromophenol	0.182	0.194	-6.6	113	0.00
15 S	2-Fluorobiphenyl	1.725	1.951	-13.1	115	-0.02
16	Acenaphthylene	2.282	2.168	5.0	98	0.00
17	Acenaphthene	1.418	1.381	2.6	100	0.00
18	Fluorene	1.782	1.694	4.9	97	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
20	4-Bromophenyl-phenylether	0.240	0.238	0.8	94	-0.01
21	Hexachlorobenzene	0.248	0.247	0.4	96	0.00
22	Pentachlorophenol	0.069	0.066	4.3	114	0.00
23	Phenanthrene	1.228	1.226	0.2	97	-0.01
24	Anthracene	1.147	1.102	3.9	95	-0.01
25 SURR	Fluoranthene-d10	5.268	4.973	5.6	93	0.00
26	Fluoranthene	1.550	1.467	5.4	95	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	98	0.00
28	Pvrene	1.680	1.661	1.1	102	0.00
29 S	Terphenyl-d14	0.852	0.931	-9.3	105	0.00
30	Benzo(a)anthracene	1.333	1.310	1.7	96	0.00
31	Chrysene	1.314	1.302	0.9	96	0.00
32	Bis(2-ethylhexyl)phthalate	3.072	2.527	17.7	85	0.00
33	Indeno(1,2,3-cd)pyrene	1.269	1.317	-3.8	105	0.00
34 I	Perylene-d12	1.000	1.000	0.0	99	0.00
35	Benzo(b)fluoranthene	1.364	1.396	-2.3	100	0.00
36	Benzo(k)fluoranthene	1.402	1.362	2.9	96	0.00
37 C	Benzo(a)pyrene	1.269	1.261	0.6	99	0.00
38	Dibenzo(a,h)anthracene	1.111	1.176	-5.9	108	0.00
39	Benzo(a,h,i)perylene	1.193	1.201	-0.7	99	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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