

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE011018\
 Data File : BE095065.D
 Acq On : 10 Jan 2018 10:11
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Jan 10 13:02:57 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 10 13:00:41 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	91	-0.01
2	1,4-Dioxane	0.394	0.383	2.8	83	0.00
3	n-Nitrosodimethylamine	0.466	0.479	-2.8	93	0.00
4 S	2-Fluorophenol	0.621	0.599	3.5	91	-0.01
5 S	Phenol-d6	0.923	0.872	5.5	89	-0.01
6	bis(2-Chloroethyl)ether	0.901	0.908	-0.8	93	-0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	98	-0.02
8 S	Nitrobenzene-d5	0.328	0.317	3.4	98	-0.02
9	Naphthalene	4.697	4.466	4.9	93	-0.02
10	Hexachlorobutadiene	0.214	0.199	7.0	90	-0.02
11 SURR	2-Methylnaphthalene-d10	0.655	0.619	5.5	93	-0.02
12	2-Methylnaphthalene	1.172	0.739	36.9#	62	-0.02
13 I	Acenaphthene-d10	1.000	1.000	0.0	90	-0.02
14 S	2,4,6-Tribromophenol	0.093	0.076	18.3	79	-0.03
15 S	2-Fluorobiphenyl	1.772	1.723	2.8	86	-0.02
16	Acenaphthylene	2.138	1.995	6.7	88	-0.03
17	Acenaphthene	1.418	1.363	3.9	87	-0.02
18	Fluorene	1.762	1.671	5.2	88	-0.03
19 I	Phenanthrene-d10	1.000	1.000	0.0	91	-0.03
20	4-Bromophenyl-phenylether	0.234	0.213	9.0	85	-0.03
21	Hexachlorobenzene	0.239	0.228	4.6	88	-0.03
22	Pentachlorophenol	0.046	0.049	-6.5	101	-0.02
23	Phenanthrene	1.246	1.252	-0.5	94	-0.03
24	Anthracene	1.279	1.239	3.1	93	-0.03
25 SURR	Fluoranthene-d10	5.189	5.356	-3.2	98	-0.03
26	Fluoranthene	1.538	1.597	-3.8	97	-0.03
27 I	Chrysene-d12	1.000	1.000	0.0	98	-0.03
28	Pyrene	1.355	1.499	-10.6	105	-0.03
29 S	Terphenyl-d14	0.753	0.757	-0.5	98	-0.03
30	Benzo(a)anthracene	1.169	1.194	-2.1	101	-0.03
31	Chrysene	1.316	1.403	-6.6	104	-0.05
32	Bis(2-ethylhexyl)phthalate	1.489	1.682	-13.0	124	-0.05
33	Indeno(1,2,3-cd)pyrene	1.071	1.060	1.0	95	-0.09
34 I	Perylene-d12	1.000	1.000	0.0	111	-0.06
35	Benzo(b)fluoranthene	1.469	1.341	8.7	101	-0.05
36	Benzo(k)fluoranthene	1.488	1.415	4.9	111	-0.06
37 C	Benzo(a)pyrene	1.303	1.247	4.3	107	-0.06
38	Dibenzo(a,h)anthracene	1.209	0.910	24.7	84	-0.08
39	Benzo(g,h,i)perylene	1.203	1.089	9.5	102	-0.09

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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