

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\BE120517\
 Data File : BE094948.D
 Acq On : 6 Dec 2017 12:43
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 06 16:48:15 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 01 13:33:28 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	84	-0.01
2	1,4-Dioxane	0.419	0.408	2.6	77	0.00
3	n-Nitrosodimethylamine	0.533	0.488	8.4	79	-0.01
4 S	2-Fluorophenol	0.633	0.646	-2.1	85	0.00
5 S	Phenol-d6	0.938	0.950	-1.3	85	0.00
6	bis(2-Chloroethyl)ether	0.930	0.855	8.1	76	-0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	81	0.00
8 S	Nitrobenzene-d5	0.237	0.310	-30.8#	107	-0.02
9	Naphthalene	4.690	4.613	1.6	79	0.00
10	Hexachlorobutadiene	0.198	0.200	-1.0	81	0.00
11 SURR	2-Methylnaphthalene-d10	0.638	0.648	-1.6	82	0.00
12	2-Methylnaphthalene	1.165	1.142	2.0	80	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	85	-0.02
14 S	2,4,6-Tribromophenol	0.086	0.112	-30.2#	114	0.00
15 S	2-Fluorobiphenyl	1.713	1.700	0.8	83	0.00
16	Acenaphthylene	2.260	2.139	5.4	82	0.00
17	Acenaphthene	1.461	1.416	3.1	84	-0.02
18	Fluorene	1.859	1.807	2.8	81	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	85	0.00
20	4-Bromophenyl-phenylether	0.223	0.221	0.9	85	0.00
21	Hexachlorobenzene	0.218	0.221	-1.4	85	0.00
22	Pentachlorophenol	0.054	0.053	1.9	92	-0.01
23	Phenanthrene	1.275	1.219	4.4	81	0.00
24	Anthracene	1.351	1.268	6.1	82	0.00
25 SURR	Fluoranthene-d10	5.466	4.972	9.0	78	0.00
26	Fluoranthene	1.638	1.492	8.9	78	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	76	-0.01
28	Pvrene	1.381	1.367	1.0	76	-0.01
29 S	Terphenyl-d14	0.739	0.742	-0.4	78	-0.01
30	Benzo(a)anthracene	1.227	1.182	3.7	76	-0.01
31	Chrysene	1.355	1.276	5.8	72	0.00
32	Bis(2-ethylhexyl)phthalate	2.033	1.886	7.2	83	-0.01
33	Indeno(1,2,3-cd)pyrene	1.109	1.048	5.5	73	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	74	-0.01
35	Benzo(b)fluoranthene	1.460	1.429	2.1	74	-0.01
36	Benzo(k)fluoranthene	1.489	1.366	8.3	70	-0.01
37 C	Benzo(a)pyrene	1.280	1.236	3.4	73	-0.01
38	Dibenzo(a,h)anthracene	1.183	1.152	2.6	74	-0.03
39	Benzo(a,h,i)perylene	1.181	1.143	3.2	73	-0.02

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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