

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\BE120117\
 Data File : BE094906.D
 Acq On : 1 Dec 2017 15:50
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 02 00:46:39 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	0.390	6.9	74	0.00
3	n-Nitrosodimethylamine	0.533	0.512	3.9	82	-0.01
4 S	2-Fluorophenol	0.633	0.608	3.9	80	0.00
5 S	Phenol-d6	0.938	0.897	4.4	80	0.00
6	bis(2-Chloroethyl)ether	0.930	0.882	5.2	78	-0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	82	0.00
8 S	Nitrobenzene-d5	0.237	0.268	-13.1	94	0.00
9	Naphthalene	4.690	4.579	2.4	80	0.00
10	Hexachlorobutadiene	0.198	0.199	-0.5	82	0.00
11 SURR	2-Methylnaphthalene-d10	0.638	0.631	1.1	81	0.00
12	2-Methylnaphthalene	1.165	1.131	2.9	81	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	84	-0.01
14 S	2,4,6-Tribromophenol	0.086	0.085	1.2	87	0.00
15 S	2-Fluorobiphenyl	1.713	1.703	0.6	82	0.00
16	Acenaphthylene	2.260	2.132	5.7	82	0.00
17	Acenaphthene	1.461	1.424	2.5	84	-0.01
18	Fluorene	1.859	1.805	2.9	81	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	84	0.00
20	4-Bromophenyl-phenylether	0.223	0.218	2.2	83	0.00
21	Hexachlorobenzene	0.218	0.231	-6.0	88	0.00
22	Pentachlorophenol	0.054	0.053	1.9	91	-0.01
23	Phenanthrene	1.275	1.253	1.7	82	0.00
24	Anthracene	1.351	1.275	5.6	82	0.00
25 SURR	Fluoranthene-d10	5.466	5.281	3.4	82	0.00
26	Fluoranthene	1.638	1.584	3.3	82	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	82	0.00
28	Pvrene	1.381	1.504	-8.9	89	-0.02
29 S	Terphenyl-d14	0.739	0.728	1.5	81	0.00
30	Benzo(a)anthracene	1.227	1.211	1.3	83	-0.02
31	Chrysene	1.355	1.346	0.7	81	0.00
32	Bis(2-ethylhexyl)phthalate	2.033	1.681	17.3	79	0.00
33	Indeno(1,2,3-cd)pyrene	1.109	1.085	2.2	81	0.00
34 I	Perylene-d12	1.000	1.000	0.0	84	0.00
35	Benzo(b)fluoranthene	1.460	1.425	2.4	84	0.00
36	Benzo(k)fluoranthene	1.489	1.431	3.9	83	0.00
37 C	Benzo(a)pyrene	1.280	1.231	3.8	83	0.00
38	Dibenzo(a,h)anthracene	1.183	1.123	5.1	82	0.00
39	Benzo(a,h,i)perylene	1.181	1.120	5.2	81	-0.01

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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