

Data Path : Z:\HPCHEM1\BNA_E\Data\BE041218\
 Data File : BE095984.D
 Acq On : 12 Apr 2018 9:51
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 12 13:55:19 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE041118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 12 13:55:20 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	105	0.00
2	1,4-Dioxane	0.618	0.584	5.5	107	0.00
3	n-Nitrosodimethylamine	0.757	0.661	12.7	99	0.00
4 S	2-Fluorophenol	1.237	1.142	7.7	102	0.00
5 S	Phenol-d6	1.707	1.561	8.6	101	0.00
6	bis(2-Chloroethyl)ether	1.502	1.417	5.7	103	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
8 S	Nitrobenzene-d5	0.469	0.433	7.7	102	0.00
9	Naphthalene	4.217	3.988	5.4	103	0.00
10	Hexachlorobutadiene	0.187	0.196	-4.8	118	0.00
11 SURR	2-Methylnaphthalene-d10	0.634	0.619	2.4	107	0.00
12	2-Methylnaphthalene	0.698	0.656	6.0	103	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
14 S	2,4,6-Tribromophenol	0.193	0.172	10.9	102	0.00
15 S	2-Fluorobiphenyl	1.711	1.621	5.3	104	0.00
16	Acenaphthylene	2.136	1.985	7.1	105	-0.01
17	Acenaphthene	1.284	1.203	6.3	106	0.00
18	Fluorene	1.588	1.499	5.6	106	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	104	-0.01
20	4-Bromophenyl-phenylether	0.238	0.229	3.8	107	0.00
21	Hexachlorobenzene	0.237	0.223	5.9	102	0.00
22	Pentachlorophenol	0.078	0.073	6.4	121	0.00
23	Phenanthrene	1.124	1.055	6.1	104	0.00
24	Anthracene	1.074	1.035	3.6	109	0.00
25 SURR	Fluoranthene-d10	5.430	5.144	5.3	107	0.00
26	Fluoranthene	1.477	1.380	6.6	105	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	104	0.00
28	Pyrene	0.788	0.718	8.9	105	0.00
29 S	Terphenyl-d14	0.884	0.836	5.4	107	0.00
30	Benzo(a)anthracene	1.283	1.233	3.9	109	0.00
31	Chrysene	1.239	1.150	7.2	101	0.00
32	Bis(2-ethylhexyl)phthalate	3.311	3.151	4.8	106	0.00
33	Indeno(1,2,3-cd)pyrene	1.201	1.133	5.7	104	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	100	0.00
35	Benzo(b)fluoranthene	1.200	1.165	2.9	102	0.00
36	Benzo(k)fluoranthene	1.234	1.163	5.8	101	0.00
37 C	Benzo(a)pyrene	1.153	1.099	4.7	102	0.00
38	Dibenzo(a,h)anthracene	1.017	0.991	2.6	106	0.00
39	Benzo(g,h,i)perylene	1.076	1.039	3.4	102	0.00

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