

Data Path : Z:\HPCHEM1\BNA E\Data\BE041118\
 Data File : BE095982.D
 Acq On : 11 Apr 2018 21:56
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 12 04:17:30 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE041118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 11 17:26:43 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	110	0.00
2	1,4-Dioxane	0.618	0.579	6.3	111	0.00
3	n-Nitrosodimethylamine	0.757	0.687	9.2	108	0.00
4 S	2-Fluorophenol	1.237	1.192	3.6	111	0.00
5 S	Phenol-d6	1.707	1.609	5.7	109	0.00
6	bis(2-Chloroethyl)ether	1.502	1.377	8.3	105	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
8 S	Nitrobenzene-d5	0.469	0.441	6.0	109	0.00
9	Naphthalene	4.217	4.011	4.9	108	0.00
10	Hexachlorobutadiene	0.187	0.193	-3.2	121	0.00
11 SURR	2-Methylnaphthalene-d10	0.634	0.606	4.4	109	0.00
12	2-Methylnaphthalene	0.698	0.672	3.7	110	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
14 S	2,4,6-Tribromophenol	0.193	0.189	2.1	119	0.00
15 S	2-Fluorobiphenyl	1.711	1.613	5.7	109	0.00
16	Acenaphthylene	2.136	1.976	7.5	110	0.00
17	Acenaphthene	1.284	1.202	6.4	112	0.00
18	Fluorene	1.588	1.486	6.4	111	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	112	0.00
20	4-Bromophenyl-phenylether	0.238	0.224	5.9	112	0.00
21	Hexachlorobenzene	0.237	0.223	5.9	109	0.00
22	Pentachlorophenol	0.078	0.073	6.4	129	0.00
23	Phenanthrene	1.124	1.041	7.4	109	0.00
24	Anthracene	1.074	1.004	6.5	113	0.00
25 SURR	Fluoranthene-d10	5.430	5.101	6.1	113	0.00
26	Fluoranthene	1.477	1.366	7.5	111	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	109	0.00
28	Pvrene	0.788	0.739	6.2	113	0.00
29 S	Terphenyl-d14	0.884	0.857	3.1	115	0.00
30	Benzo(a)anthracene	1.283	1.192	7.1	110	0.00
31	Chrysene	1.239	1.163	6.1	108	0.00
32	Bis(2-ethylhexyl)phthalate	3.311	3.223	2.7	114	0.00
33	Indeno(1,2,3-cd)pyrene	1.201	1.141	5.0	110	0.00
34 I	Perylene-d12	1.000	1.000	0.0	107	0.00
35	Benzo(b)fluoranthene	1.200	1.129	5.9	106	0.00
36	Benzo(k)fluoranthene	1.234	1.146	7.1	107	0.00
37 C	Benzo(a)pyrene	1.153	1.076	6.7	107	0.00
38	Dibenzo(a,h)anthracene	1.017	0.987	2.9	113	0.00
39	Benzo(a,h,i)perylene	1.076	0.997	7.3	105	0.00

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