

Data Path : Z:\HPCHEM1\BNA E\Data\BE043018\
 Data File : BE096213.D
 Acq On : 30 Apr 2018 9:58
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 30 14:27:35 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE041118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 30 14:27:29 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	53	0.00
2	1,4-Dioxane	0.618	0.595	3.7	54	0.00
3	n-Nitrosodimethylamine	0.757	0.828	-9.4	63	0.00
4 S	2-Fluorophenol	1.237	1.183	4.4	53	0.00
5 S	Phenol-d6	1.707	1.623	4.9	53	0.00
6	bis(2-Chloroethyl)ether	1.502	1.456	3.1	53	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	52	0.00
8 S	Nitrobenzene-d5	0.469	0.457	2.6	55	0.00
9	Naphthalene	4.217	3.938	6.6	51	0.00
10	Hexachlorobutadiene	0.187	0.116	38.0#	35#	0.00
11 SURR	2-Methylnaphthalene-d10	0.634	0.580	8.5	51	0.00
12	2-Methylnaphthalene	0.698	0.650	6.9	52	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	52	0.00
14 S	2,4,6-Tribromophenol	0.193	0.172	10.9	50	0.00
15 S	2-Fluorobiphenyl	1.711	1.654	3.3	52	0.00
16	Acenaphthylene	2.136	2.042	4.4	53	0.00
17	Acenaphthene	1.284	1.199	6.6	52	0.00
18	Fluorene	1.588	1.501	5.5	52	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	53	0.00
20	4-Bromophenyl-phenylether	0.238	0.219	8.0	52	0.00
21	Hexachlorobenzene	0.237	0.233	1.7	54	0.00
22	Pentachlorophenol	0.078	0.065	16.7	55	0.00
23	Phenanthrene	1.124	1.085	3.5	55	0.00
24	Anthracene	1.074	1.015	5.5	55	0.00
25 SURR	Fluoranthene-d10	5.430	5.107	5.9	54	0.00
26	Fluoranthene	1.477	1.428	3.3	56	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	52	0.00
28	Pvrene	0.788	0.047	94.0#	3#	0.00
29 S	Terphenyl-d14	0.884	0.841	4.9	54	0.00
30	Benzo(a)anthracene	1.283	1.239	3.4	55	0.00
31	Chrysene	1.239	1.237	0.2	55	0.00
32	Bis(2-ethylhexyl)phthalate	3.311	3.241	2.1	55	0.00
33	Indeno(1,2,3-cd)pyrene	1.201	1.197	0.3	55	0.00
34 I	Perylene-d12	1.000	1.000	0.0	50	0.00
35	Benzo(b)fluoranthene	1.200	1.206	-0.5	53	0.00
36	Benzo(k)fluoranthene	1.234	1.194	3.2	52	0.00
37 C	Benzo(a)pyrene	1.153	1.125	2.4	52	0.00
38	Dibenzo(a,h)anthracene	1.017	1.052	-3.4	56	0.00
39	Benzo(a,h,i)perylene	1.076	1.071	0.5	53	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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