

Data Path : Z:\HPCHEM1\BNA E\Data\BE043018\
 Data File : BE096229.D
 Acq On : 30 Apr 2018 19:35
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: May 01 05:14:00 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	58	0.00
2	1,4-Dioxane	0.400	0.346	13.5	54	0.00
3	n-Nitrosodimethylamine	0.400	0.396	1.0	62	0.00
4 S	2-Fluorophenol	0.400	0.387	3.3	59	0.01
5 S	Phenol-d6	0.400	0.397	0.8	60	0.01
6	bis(2-Chloroethyl)ether	0.400	0.365	8.8	55	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	55	0.00
8 S	Nitrobenzene-d5	0.400	0.400	0.0	60	0.00
9	Naphthalene	0.400	0.373	6.8	55	0.00
10	Hexachlorobutadiene	0.400	0.246	38.5#	38	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.367	8.3	54	0.00
12	2-Methylnaphthalene	0.400	0.365	8.8	55	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	57	0.00
14 S	2,4,6-Tribromophenol	0.400	0.375	6.3	59	0.00
15 S	2-Fluorobiphenyl	0.400	0.367	8.3	55	0.00
16	Acenaphthylene	0.400	0.372	7.0	57	0.00
17	Acenaphthene	0.400	0.366	8.5	56	0.00
18	Fluorene	0.400	0.366	8.5	56	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	58	0.00
20	4-Bromophenyl-phenylether	0.400	0.366	8.5	56	0.00
21	Hexachlorobenzene	0.400	0.361	9.8	54	0.00
22	Pentachlorophenol	0.400	0.405	-1.3	67	0.00
23	Phenanthrene	0.400	0.385	3.8	59	0.00
24	Anthracene	0.400	0.379	5.3	59	0.00
25 SURR	Fluoranthene-d10	0.400	0.373	6.8	58	0.00
26	Fluoranthene	0.400	0.386	3.5	60	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	53	0.01
28	Pvrene	0.400	0.032	92.0#	5	0.00
29 S	Terphenyl-d14	0.400	0.406	-1.5	58	0.00
30	Benzo(a)anthracene	0.400	0.388	3.0	56	0.00
31	Chrysene	0.400	0.402	-0.5	56	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.429	-7.2	61	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.414	-3.5	58	0.00
34 I	Perylene-d12	0.400	0.400	0.0	53	0.00
35	Benzo(b)fluoranthene	0.400	0.392	2.0	55	0.00
36	Benzo(k)fluoranthene	0.400	0.388	3.0	55	0.00
37 C	Benzo(a)pyrene	0.400	0.388	3.0	55	0.00
38	Dibenzo(a,h)anthracene	0.400	0.413	-3.2	59	0.00
39	Benzo(a,h,i)perylene	0.400	0.397	0.8	56	0.00

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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