

Data Path : Z:\HPCHEM1\BNA_E\Data\BE112717\
 Data File : BE094863.D
 Acq On : 27 Nov 2017 9:23
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 27 12:13:12 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 23 22:40:47 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	48	0.01
2	1,4-Dioxane	0.400	0.397	0.8	41	-0.04
3	n-Nitrosodimethylamine	0.400	0.308	23.0	37	0.00
4 S	2-Fluorophenol	0.400	0.336	16.0	40	0.09
5 S	Phenol-d6	0.400	0.387	3.3	48	0.08
6	bis(2-Chloroethyl)ether	0.400	0.306	23.5	37	0.04
7 I	Naphthalene-d8	0.400	0.400	0.0	58	0.00
8 S	Nitrobenzene-d5	0.400	0.317	20.8	48	0.07
9	Naphthalene	0.400	0.338	15.5	48	0.02
10	Hexachlorobutadiene	0.400	0.393	1.8	56	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.382	4.5	55	0.02
12	2-Methylnaphthalene	0.400	0.366	8.5	53	0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	67	0.00
14 S	2,4,6-Tribromophenol	0.400	0.372	7.0	64	0.03
15 S	2-Fluorobiphenyl	0.400	0.339	15.3	56	0.01
16	Acenaphthylene	0.400	0.337	15.8	57	0.00
17	Acenaphthene	0.400	0.363	9.3	59	-0.01
18	Fluorene	0.400	0.375	6.3	62	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	72	0.03
20	4-Bromophenyl-phenylether	0.400	0.400	0.0	74	0.00
21	Hexachlorobenzene	0.400	0.546	-36.5#	86	0.00
22	Pentachlorophenol	0.400	0.459	-14.8	83	0.03
23	Phenanthrene	0.400	0.468	-17.0	77	0.00
24	Anthracene	0.400	0.431	-7.7	76	0.00
25 SURR	Fluoranthene-d10	0.400	0.464	-16.0	75	0.00
26	Fluoranthene	0.400	0.452	-13.0	73	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	78	0.00
28	Pyrene	0.400	0.368	8.0	74	0.00
29 S	Terphenyl-d14	0.400	0.365	8.8	73	0.00
30	Benzo(a)anthracene	0.400	0.370	7.5	73	0.01
31	Chrysene	0.400	0.400	0.0	81	0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.332	17.0	71	-0.02
33	Indeno(1,2,3-cd)pyrene	0.400	0.406	-1.5	81	0.05
34 I	Perylene-d12	0.400	0.400	0.0	79	0.02
35	Benzo(b)fluoranthene	0.400	0.386	3.5	78	0.01
36	Benzo(k)fluoranthene	0.400	0.382	4.5	76	0.02
37 C	Benzo(a)pyrene	0.400	0.386	3.5	77	0.02
38	Dibenzo(a,h)anthracene	0.400	0.396	1.0	80	0.03
39	Benzo(g,h,i)perylene	0.400	0.433	-8.2	87	0.06

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

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