

Data Path : Z:\HPCHEM1\BNA_E\Data\BE070117\
 Data File : BE093386.D
 Acq On : 1 Jul 2017 11:08
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 03 05:21:10 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 01 15:43:24 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	93	-0.01
2	1,4-Dioxane	0.400	0.387	3.3	93	0.00
3	n-Nitrosodimethylamine	0.400	0.386	3.5	91	-0.01
4 S	2-Fluorophenol	0.400	0.386	3.5	93	0.00
5 S	Phenol-d6	0.400	0.366	8.5	91	0.00
6	bis(2-Chloroethyl)ether	0.400	0.396	1.0	100	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	92	0.00
8 S	Nitrobenzene-d5	0.400	0.384	4.0	102	-0.02
9	Naphthalene	0.400	0.393	1.8	94	-0.02
10	Hexachlorobutadiene	0.400	0.408	-2.0	95	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.383	4.3	92	0.00
12	2-Methylnaphthalene	0.400	0.382	4.5	91	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	88	-0.02
14 S	2,4,6-Tribromophenol	0.400	0.379	5.3	84	0.00
15 S	2-Fluorobiphenyl	0.400	0.415	-3.7	92	0.00
16	Acenaphthylene	0.400	0.372	7.0	86	-0.02
17	Acenaphthene	0.400	0.393	1.8	89	0.00
18	Fluorene	0.400	0.404	-1.0	90	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	85	0.00
20	4-Bromophenyl-phenylether	0.400	0.542	-35.5#	116	0.00
21	Hexachlorobenzene	0.400	0.478	-19.5	101	0.00
22	Pentachlorophenol	0.400	0.421	-5.2	92	-0.03
23	Phenanthrene	0.400	0.455	-13.7	93	0.00
24	Anthracene	0.400	0.339	15.3	78	0.00
25 SURR	Fluoranthene-d10	0.400	0.397	0.8	89	0.00
26	Fluoranthene	0.400	0.396	1.0	89	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	94	-0.01
28	Pyrene	0.400	0.375	6.3	90	0.00
29 S	Terphenyl-d14	0.400	0.388	3.0	93	0.00
30	Benzo(a)anthracene	0.400	0.378	5.5	93	0.00
31	Chrysene	0.400	0.395	1.3	94	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.327	18.3	92	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.403	-0.8	98	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	97	0.00
35	Benzo(b)fluoranthene	0.400	0.387	3.3	98	0.00
36	Benzo(k)fluoranthene	0.400	0.386	3.5	98	-0.01
37 C	Benzo(a)pyrene	0.400	0.380	5.0	97	0.00
38	Dibenzo(a,h)anthracene	0.400	0.386	3.5	98	0.00
39	Benzo(g,h,i)perylene	0.400	0.390	2.5	98	0.00

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