

Data Path : Z:\HPCHEM1\BNA E\DATA\BE051617\
 Data File : BE093051.D
 Acq On : 16 May 2017 13:27
 Operator : SJ/MA
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: May 18 00:54:27 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE051517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 15 13:35:57 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	128	0.00
2	1,4-Dioxane	0.400	0.369	7.8	136	0.00
3	n-Nitrosodimethylamine	0.400	0.454	-13.5	167	-0.01
4 S	2-Fluorophenol	0.400	0.378	5.5	137	0.00
5 S	Phenol-d6	0.400	0.438	-9.5	160	0.00
6	bis(2-Chloroethyl)ether	0.400	0.404	-1.0	148	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	139	0.00
8 S	Nitrobenzene-d5	0.400	0.448	-12.0	181	0.00
9	Naphthalene	0.400	0.394	1.5	153	-0.02
10	Hexachlorobutadiene	0.400	0.381	4.8	150	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.361	9.8	141	0.00
12	2-Methylnaphthalene	0.400	0.353	11.8	139	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	134	0.00
14 S	2,4,6-Tribromophenol	0.400	0.544	-36.0#	212	-0.01
15 S	2-Fluorobiphenyl	0.400	0.380	5.0	142	0.00
16	Acenaphthylene	0.400	0.412	-3.0	163	0.00
17	Acenaphthene	0.400	0.384	4.0	147	0.00
18	Fluorene	0.400	0.398	0.5	151	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	139	0.00
20	4-Bromophenyl-phenylether	0.400	0.404	-1.0	176	0.00
21	Hexachlorobenzene	0.400	0.257	35.8#	137	0.00
22	Pentachlorophenol	0.400	0.597	-49.2#	290	-0.03
23	Phenanthrene	0.400	0.352	12.0	156	0.00
24	Anthracene	0.400	0.317	20.8	160	0.00
25 SURR	Fluoranthene-d10	0.400	0.385	3.8	164	0.00
26	Fluoranthene	0.400	0.362	9.5	159	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	152	-0.02
28	Pvrene	0.400	0.379	5.3	157	0.00
29 S	Terphenyl-d14	0.400	0.380	5.0	159	0.00
30	Benzo(a)anthracene	0.400	0.426	-6.5	182	0.00
31	Chrysene	0.400	0.383	4.3	164	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.726	-81.5#	365	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.415	-3.7	179	0.00
34 I	Perylene-d12	0.400	0.400	0.0	157	-0.01
35	Benzo(b)fluoranthene	0.400	0.394	1.5	181	0.00
36	Benzo(k)fluoranthene	0.400	0.363	9.3	162	0.00
37 C	Benzo(a)pyrene	0.400	0.385	3.8	177	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.410	-2.5	186	0.00
39	Benzo(a,h,i)perylene	0.400	0.384	4.0	173	-0.02

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

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