

Data Path : Z:\HPCHEM1\BNA E\DATA\BE051917\
 Data File : BE093063.D
 Acq On : 19 May 2017 16:29
 Operator : SJ/MA
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: May 19 18:06:36 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	80	0.00
2	1,4-Dioxane	0.400	0.368	8.0	85	0.00
3	n-Nitrosodimethylamine	0.400	0.358	10.5	82	0.00
4 S	2-Fluorophenol	0.400	0.356	11.0	80	0.00
5 S	Phenol-d6	0.400	0.353	11.8	81	0.00
6	bis(2-Chloroethyl)ether	0.400	0.353	11.8	81	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	82	0.00
8 S	Nitrobenzene-d5	0.400	0.372	7.0	88	0.00
9	Naphthalene	0.400	0.352	12.0	81	-0.02
10	Hexachlorobutadiene	0.400	0.370	7.5	86	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.351	12.3	81	0.00
12	2-Methylnaphthalene	0.400	0.350	12.5	81	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	82	0.00
14 S	2,4,6-Tribromophenol	0.400	0.381	4.8	91	-0.01
15 S	2-Fluorobiphenyl	0.400	0.352	12.0	80	0.00
16	Acenaphthylene	0.400	0.346	13.5	84	0.00
17	Acenaphthene	0.400	0.352	12.0	82	0.00
18	Fluorene	0.400	0.354	11.5	82	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	86	0.00
20	4-Bromophenyl-phenylether	0.400	0.357	10.8	97	0.00
21	Hexachlorobenzene	0.400	0.204	49.0#	73	0.00
22	Pentachlorophenol	0.400	0.308	23.0	97	0.00
23	Phenanthrene	0.400	0.301	24.8	86	0.00
24	Anthracene	0.400	0.247	38.3#	83	0.00
25 SURR	Fluoranthene-d10	0.400	0.318	20.5	85	0.00
26	Fluoranthene	0.400	0.324	19.0	89	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	88	-0.02
28	Pvrene	0.400	0.349	12.8	84	-0.02
29 S	Terphenyl-d14	0.400	0.317	20.8	76	0.00
30	Benzo(a)anthracene	0.400	0.374	6.5	92	0.00
31	Chrysene	0.400	0.335	16.3	83	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.356	11.0	93	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.365	8.8	91	0.00
34 I	Perylene-d12	0.400	0.400	0.0	87	0.00
35	Benzo(b)fluoranthene	0.400	0.378	5.5	96	0.00
36	Benzo(k)fluoranthene	0.400	0.320	20.0	79	0.00
37 C	Benzo(a)pyrene	0.400	0.346	13.5	88	0.00
38	Dibenzo(a,h)anthracene	0.400	0.368	8.0	93	0.00
39	Benzo(a,h,i)perylene	0.400	0.361	9.8	90	-0.02

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

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