

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE011018\
 Data File : BE095065.D
 Acq On : 10 Jan 2018 10:11
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Jan 10 13:02:57 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 10 13:00:41 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	91	-0.01
2	1,4-Dioxane	0.400	0.388	3.0	83	0.00
3	n-Nitrosodimethylamine	0.400	0.411	-2.7	93	0.00
4 S	2-Fluorophenol	0.400	0.386	3.5	91	-0.01
5 S	Phenol-d6	0.400	0.378	5.5	89	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.403	-0.8	93	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	98	-0.02
8 S	Nitrobenzene-d5	0.400	0.386	3.5	98	-0.02
9	Naphthalene	0.400	0.380	5.0	93	-0.02
10	Hexachlorobutadiene	0.400	0.372	7.0	90	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.378	5.5	93	-0.02
12	2-Methylnaphthalene	0.400	0.252	37.0#	62	-0.02
13 I	Acenaphthene-d10	0.400	0.400	0.0	90	-0.02
14 S	2,4,6-Tribromophenol	0.400	0.328	18.0	79	-0.03
15 S	2-Fluorobiphenyl	0.400	0.389	2.8	86	-0.02
16	Acenaphthylene	0.400	0.373	6.8	88	-0.03
17	Acenaphthene	0.400	0.384	4.0	87	-0.02
18	Fluorene	0.400	0.379	5.3	88	-0.03
19 I	Phenanthrene-d10	0.400	0.400	0.0	91	-0.03
20	4-Bromophenyl-phenylether	0.400	0.363	9.3	85	-0.03
21	Hexachlorobenzene	0.400	0.381	4.8	88	-0.03
22	Pentachlorophenol	0.400	0.410	-2.5	101	-0.02
23	Phenanthrene	0.400	0.402	-0.5	94	-0.03
24	Anthracene	0.400	0.388	3.0	93	-0.03
25 SURR	Fluoranthene-d10	0.400	0.413	-3.2	98	-0.03
26	Fluoranthene	0.400	0.415	-3.7	97	-0.03
27 I	Chrysene-d12	0.400	0.400	0.0	98	-0.03
28	Pyrene	0.400	0.443	-10.7	105	-0.03
29 S	Terphenyl-d14	0.400	0.402	-0.5	98	-0.03
30	Benzo(a)anthracene	0.400	0.409	-2.2	101	-0.03
31	Chrysene	0.400	0.426	-6.5	104	-0.05
32	Bis(2-ethylhexyl)phthalate	0.400	0.452	-13.0	124	-0.05
33	Indeno(1,2,3-cd)pyrene	0.400	0.396	1.0	95	-0.09
34 I	Perylene-d12	0.400	0.400	0.0	111	-0.06
35	Benzo(b)fluoranthene	0.400	0.365	8.8	101	-0.05
36	Benzo(k)fluoranthene	0.400	0.380	5.0	111	-0.06
37 C	Benzo(a)pyrene	0.400	0.383	4.3	107	-0.06
38	Dibenzo(a,h)anthracene	0.400	0.301	24.8	84	-0.08
39	Benzo(g,h,i)perylene	0.400	0.362	9.5	102	-0.09

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

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