

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE092618\
 Data File : BE097688.D
 Acq On : 26 Sep 2018 17:15
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 26 18:01:43 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE092418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 24 18:40:34 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	98	-0.01
2	1,4-Dioxane	0.400	0.371	7.3	99	0.00
3	n-Nitrosodimethylamine	0.400	0.490	-22.5	141	-0.01
4 S	2-Fluorophenol	0.400	0.469	-17.2	130	-0.01
5 S	Phenol-d6	0.400	0.490	-22.5	137	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.450	-12.5	111	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	104	-0.01
8 S	Nitrobenzene-d5	0.400	0.447	-11.7	133	-0.01
9	Naphthalene	0.400	0.409	-2.2	114	-0.01
10	Hexachlorobutadiene	0.400	0.390	2.5	110	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.421	-5.2	117	0.00
12	2-Methylnaphthalene	0.400	0.442	-10.5	123	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	107	0.00
14 S	2,4,6-Tribromophenol	0.400	0.419	-4.7	135	0.00
15 S	2-Fluorobiphenyl	0.400	0.413	-3.2	118	-0.01
16	Acenaphthylene	0.400	0.408	-2.0	122	0.00
17	Acenaphthene	0.400	0.420	-5.0	121	0.00
18	Fluorene	0.400	0.404	-1.0	118	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	99	0.00
20	4-Bromophenyl-phenylether	0.400	0.422	-5.5	112	0.00
21	Hexachlorobenzene	0.400	0.411	-2.7	109	0.00
22	Pentachlorophenol	0.400	0.539	-34.8#	160	-0.01
23	Phenanthrene	0.400	0.409	-2.2	109	0.00
24	Anthracene	0.400	0.420	-5.0	116	0.00
25 SURR	Fluoranthene-d10	0.400	0.391	2.3	107	0.00
26	Fluoranthene	0.400	0.389	2.8	107	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	96	0.00
28	Pvrene	0.400	0.421	-5.2	108	0.00
29 S	Terphenyl-d14	0.400	0.410	-2.5	107	0.00
30	Benzo(a)anthracene	0.400	0.412	-3.0	113	0.00
31	Chrysene	0.400	0.411	-2.7	101	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.536	-34.0#	146	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.421	-5.2	107	0.00
34 I	Perylene-d12	0.400	0.400	0.0	97	0.00
35	Benzo(b)fluoranthene	0.400	0.410	-2.5	108	0.00
36	Benzo(k)fluoranthene	0.400	0.402	-0.5	105	0.00
37 C	Benzo(a)pyrene	0.400	0.409	-2.2	110	0.00
38	Dibenzo(a,h)anthracene	0.400	0.406	-1.5	111	0.00
39	Benzo(a,h,i)perylene	0.400	0.405	-1.3	107	0.00

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

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