

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121118\
 Data File : BE098352.D
 Acq On : 12 Dec 2018 3:33
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 12 04:09:26 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 10 14:59:55 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	117	0.00
2	1,4-Dioxane	0.400	0.412	-3.0	123	0.00
3	n-Nitrosodimethylamine	0.400	0.299	25.3#	95	0.00
4 S	2-Fluorophenol	0.400	0.399	0.3	121	0.00
5 S	Phenol-d6	0.400	0.377	5.8	110	0.00
6	bis(2-Chloroethyl)ether	0.400	0.375	6.3	114	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	113	-0.01
8 S	Nitrobenzene-d5	0.400	0.415	-3.7	127	0.00
9	Naphthalene	0.400	0.399	0.3	115	0.00
10	Hexachlorobutadiene	0.400	0.397	0.8	115	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.383	4.3	109	0.00
12	2-Methylnaphthalene	0.400	0.374	6.5	108	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	109	0.00
14 S	2,4,6-Tribromophenol	0.400	0.416	-4.0	120	0.00
15 S	2-Fluorobiphenyl	0.400	0.374	6.5	112	0.00
16	Acenaphthylene	0.400	0.392	2.0	111	0.00
17	Acenaphthene	0.400	0.388	3.0	107	0.00
18	Fluorene	0.400	0.395	1.3	109	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	104	-0.01
20	4-Bromophenyl-phenylether	0.400	0.369	7.8	100	0.00
21	Hexachlorobenzene	0.400	0.387	3.3	103	0.00
22	Pentachlorophenol	0.400	0.378	5.5	100	0.00
23	Phenanthrene	0.400	0.391	2.3	103	0.00
24	Anthracene	0.400	0.359	10.3	95	0.00
25 SURR	Fluoranthene-d10	0.400	0.396	1.0	100	0.00
26	Fluoranthene	0.400	0.401	-0.3	100	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	123	0.00
28	Pvrene	0.400	0.337	15.8	102	0.00
29 S	Terphenyl-d14	0.400	0.328	18.0	102	0.00
30	Benzo(a)anthracene	0.400	0.369	7.8	116	-0.01
31	Chrysene	0.400	0.379	5.3	118	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.342	14.5	106	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.393	1.8	128	0.00
34 I	Perylene-d12	0.400	0.400	0.0	130	0.00
35	Benzo(b)fluoranthene	0.400	0.385	3.8	129	0.00
36	Benzo(k)fluoranthene	0.400	0.410	-2.5	139	0.00
37 C	Benzo(a)pyrene	0.400	0.387	3.3	128	0.00
38	Dibenzo(a,h)anthracene	0.400	0.384	4.0	129	0.00
39	Benzo(a,h,i)perylene	0.400	0.379	5.3	127	0.00

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

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