

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE041519\
 Data File : BE099328.D
 Acq On : 15 Apr 2019 16:29
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 15 19:03:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE040219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 02 17:16:42 2019
 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	161	-0.03
2	1,4-Dioxane	0.400	0.426	-6.5	172	-0.01
3	n-Nitrosodimethylamine	0.400	0.405	-1.3	169	-0.03
4 S	2-Fluorophenol	0.400	0.407	-1.7	157	-0.03
5 S	Phenol-d6	0.400	0.435	-8.7	173	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.425	-6.2	171	-0.03
7 I	Naphthalene-d8	0.400	0.400	0.0	163	-0.01
8 S	Nitrobenzene-d5	0.400	0.476	-19.0	203	-0.03
9	Naphthalene	0.400	0.435	-8.7	176	-0.03
10	Hexachlorobutadiene	0.400	0.419	-4.7	170	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.430	-7.5	178	-0.02
12	2-Methylnaphthalene	0.400	0.426	-6.5	175	-0.02
13 I	Acenaphthene-d10	0.400	0.400	0.0	172	-0.03
14 S	2,4,6-Tribromophenol	0.400	0.392	2.0	188	-0.01
15 S	2-Fluorobiphenyl	0.400	0.420	-5.0	181	-0.02
16	Acenaphthylene	0.400	0.398	0.5	177	-0.02
17	Acenaphthene	0.400	0.422	-5.5	186	-0.02
18	Fluorene	0.400	0.415	-3.7	178	-0.03
19 I	Phenanthrene-d10	0.400	0.400	0.0	166	-0.03
20	4-Bromophenyl-phenylether	0.400	0.402	-0.5	170	-0.03
21	Hexachlorobenzene	0.400	0.415	-3.7	174	-0.03
22	Pentachlorophenol	0.400	0.475	-18.7	225	-0.03
23	Phenanthrene	0.400	0.434	-8.5	178	-0.03
24	Anthracene	0.400	0.409	-2.2	170	-0.01
25 SURR	Fluoranthene-d10	0.400	0.393	1.8	161	-0.02
26	Fluoranthene	0.400	0.397	0.8	163	-0.02
27 I	Chrysene-d12	0.400	0.400	0.0	134	-0.02
28	Pvrene	0.400	0.474	-18.5	162	-0.02
29 S	Terphenyl-d14	0.400	0.461	-15.3	156	-0.02
30	Benzo(a)anthracene	0.400	0.406	-1.5	136	-0.02
31	Chrysene	0.400	0.428	-7.0	142	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.361	9.8	125	-0.02
33	Indeno(1,2,3-cd)pyrene	0.400	0.399	0.3	130	-0.06
34 I	Perylene-d12	0.400	0.400	0.0	125	-0.03
35	Benzo(b)fluoranthene	0.400	0.417	-4.2	135	-0.03
36	Benzo(k)fluoranthene	0.400	0.440	-10.0	141	-0.03
37 C	Benzo(a)pyrene	0.400	0.411	-2.7	132	-0.03
38	Dibenzo(a,h)anthracene	0.400	0.413	-3.2	131	-0.06
39	Benzo(a,h,i)perylene	0.400	0.412	-3.0	132	-0.06

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

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