

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE043019\
 Data File : BE099416.D
 Acq On : 30 Apr 2019 8:57
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC00.4

Quant Time: Apr 30 09:28:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE042919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 29 18:36:43 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	91	0.00
2	1,4-Dioxane	0.550	0.547	0.5	105	0.00
3	n-Nitrosodimethylamine	0.332	0.308	7.2	95	0.00
4 S	2-Fluorophenol	1.106	1.092	1.3	94	0.00
5 S	Phenol-d6	1.471	1.484	-0.9	94	-0.01
6	bis(2-Chloroethyl)ether	1.269	1.222	3.7	89	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	94	0.00
8 S	Nitrobenzene-d5	0.375	0.369	1.6	93	0.00
9	Naphthalene	4.359	4.462	-2.4	95	0.00
10	Hexachlorobutadiene	0.303	0.310	-2.3	94	0.00
11 SURR	2-Methylnaphthalene-d10	0.721	0.725	-0.6	96	0.00
12	2-Methylnaphthalene	0.758	0.760	-0.3	95	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
14 S	2,4,6-Tribromophenol	0.252	0.224	11.1	89	0.00
15 S	2-Fluorobiphenyl	1.526	1.528	-0.1	95	-0.01
16	Acenaphthylene	1.957	1.885	3.7	93	0.00
17	Acenaphthene	1.118	1.118	0.0	97	0.00
18	Fluorene	1.649	1.622	1.6	94	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
20	4-Bromophenyl-phenylether	0.233	0.231	0.9	93	0.00
21	Hexachlorobenzene	0.222	0.224	-0.9	95	0.00
22	Pentachlorophenol	0.100	0.081	19.0	91	0.00
23	Phenanthrene	1.032	1.056	-2.3	96	0.00
24	Anthracene	0.984	0.967	1.7	94	0.00
25 SURR	Fluoranthene-d10	5.391	5.460	-1.3	96	0.00
26	Fluoranthene	1.546	1.588	-2.7	97	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	101	0.00
28	Pyrene	1.043	1.133	-8.6	96	0.00
29 S	Terphenyl-d14	0.733	0.788	-7.5	96	0.00
30	Benzo(a)anthracene	1.274	1.427	-12.0	100	-0.01
31	Chrysene	1.247	1.392	-11.6	99	0.00
32	Bis(2-ethylhexyl)phthalate	2.216	2.023	8.7	84	0.00
33	Indeno(1,2,3-cd)pyrene	1.474	1.636	-11.0	98	0.00
34 I	Perylene-d12	1.000	1.000	0.0	99	0.00
35	Benzo(b)fluoranthene	1.153	1.209	-4.9	102	0.00
36	Benzo(k)fluoranthene	1.112	1.141	-2.6	96	0.00
37 C	Benzo(a)pyrene	1.090	1.120	-2.8	98	0.00
38	Dibenzo(a,h)anthracene	1.115	1.132	-1.5	97	0.00
39	Benzo(g,h,i)perylene	1.118	1.153	-3.1	98	0.00

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