

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE022719\
 Data File : BE098781.D
 Acq On : 27 Feb 2019 17:46
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Feb 28 02:43:14 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE022619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 26 15:44:42 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	107	0.00
2	1,4-Dioxane	0.690	0.663	3.9	104	0.00
3	n-Nitrosodimethylamine	0.758	0.693	8.6	96	0.00
4 S	2-Fluorophenol	1.043	1.014	2.8	106	0.00
5 S	Phenol-d6	1.560	1.536	1.5	108	0.00
6	bis(2-Chloroethyl)ether	1.071	0.867	19.0	88	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
8 S	Nitrobenzene-d5	0.380	0.386	-1.6	107	0.00
9	Naphthalene	4.611	4.568	0.9	107	0.00
10	Hexachlorobutadiene	0.306	0.301	1.6	106	0.00
11 SURR	2-Methylnaphthalene-d10	0.757	0.726	4.1	106	0.00
12	2-Methylnaphthalene	0.735	0.726	1.2	110	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	112	0.00
14 S	2,4,6-Tribromophenol	0.158	0.180	-13.9	125	-0.01
15 S	2-Fluorobiphenyl	1.693	1.670	1.4	110	0.00
16	Acenaphthylene	2.045	1.948	4.7	111	-0.01
17	Acenaphthene	1.209	1.149	5.0	108	0.00
18	Fluorene	1.673	1.612	3.6	111	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
20	4-Bromophenyl-phenylether	0.225	0.218	3.1	117	0.00
21	Hexachlorobenzene	0.270	0.265	1.9	109	0.00
22	Pentachlorophenol	0.061	0.049	19.7	111	0.00
23	Phenanthrene	1.096	1.052	4.0	111	0.00
24	Anthracene	0.885	0.791	10.6	107	0.00
25 SURR	Fluoranthene-d10	5.692	5.341	6.2	108	0.00
26	Fluoranthene	1.414	1.349	4.6	111	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	120	-0.01
28	Pyrene	1.329	1.160	12.7	110	0.00
29 S	Terphenyl-d14	0.992	0.873	12.0	110	0.00
30	Benzo(a)anthracene	1.152	1.030	10.6	117	0.00
31	Chrysene	1.300	1.116	14.2	102	0.00
32	Bis(2-ethylhexyl)phthalate	2.456	1.888	23.1	102	0.00
33	Indeno(1,2,3-cd)pyrene	1.032	0.945	8.4	107	0.00
34 I	Perylene-d12	1.000	1.000	0.0	109	0.00
35	Benzo(b)fluoranthene	1.260	1.110	11.9	98	0.00
36	Benzo(k)fluoranthene	1.313	1.169	11.0	100	0.00
37 C	Benzo(a)pyrene	1.163	1.064	8.5	103	0.00
38	Dibenzo(a,h)anthracene	0.739	0.828	-12.0	130	0.00
39	Benzo(g,h,i)perylene	1.085	0.943	13.1	96	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0