

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE041919\
 Data File : BE099346.D
 Acq On : 19 Apr 2019 13:15
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC00.4

Quant Time: Apr 19 22:56:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 19 06:51:38 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	105	0.00
2	1,4-Dioxane	0.560	0.587	-4.8	111	0.00
3	n-Nitrosodimethylamine	0.301	0.279	7.3	115	0.00
4 S	2-Fluorophenol	0.977	1.043	-6.8	112	0.00
5 S	Phenol-d6	1.395	1.446	-3.7	109	0.00
6	bis(2-Chloroethyl)ether	1.301	1.300	0.1	101	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
8 S	Nitrobenzene-d5	0.311	0.317	-1.9	108	0.00
9	Naphthalene	4.364	4.501	-3.1	103	0.00
10	Hexachlorobutadiene	0.277	0.293	-5.8	105	0.00
11 SURR	2-Methylnaphthalene-d10	0.706	0.719	-1.8	103	0.00
12	2-Methylnaphthalene	0.743	0.776	-4.4	106	0.01
13 I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
14 S	2,4,6-Tribromophenol	0.133	0.149	-12.0	113	0.00
15 S	2-Fluorobiphenyl	1.572	1.608	-2.3	101	0.00
16	Acenaphthylene	1.808	1.792	0.9	103	0.00
17	Acenaphthene	1.114	1.154	-3.6	104	0.00
18	Fluorene	1.601	1.630	-1.8	101	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
20	4-Bromophenyl-phenylether	0.222	0.221	0.5	100	0.00
21	Hexachlorobenzene	0.213	0.223	-4.7	103	0.00
22	Pentachlorophenol	0.057	0.057	0.0	122	0.00
23	Phenanthrene	1.059	1.084	-2.4	102	0.00
24	Anthracene	0.942	0.953	-1.2	103	0.00
25 SURR	Fluoranthene-d10	4.806	4.774	0.7	102	0.00
26	Fluoranthene	1.418	1.409	0.6	102	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	103	0.00
28	Pyrene	1.178	1.184	-0.5	101	0.00
29 S	Terphenyl-d14	0.787	0.795	-1.0	102	0.00
30	Benzo(a)anthracene	1.176	1.236	-5.1	113	0.00
31	Chrysene	1.251	1.259	-0.6	99	0.00
32	Bis(2-ethylhexyl)phthalate	1.070	1.140	-6.5	128	0.00
33	Indeno(1,2,3-cd)pyrene	1.236	1.296	-4.9	108	0.00
34 I	Perylene-d12	1.000	1.000	0.0	106	0.00
35	Benzo(b)fluoranthene	1.225	1.233	-0.7	107	0.00
36	Benzo(k)fluoranthene	1.221	1.226	-0.4	104	0.00
37 C	Benzo(a)pyrene	1.097	1.116	-1.7	108	0.00
38	Dibenzo(a,h)anthracene	1.121	1.116	0.4	105	0.00
39	Benzo(g,h,i)perylene	1.137	1.144	-0.6	106	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE041919\
Data File : BE099346.D
Acq On : 19 Apr 2019 13:15
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Apr 19 22:56:58 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE041819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 19 06:51:38 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)

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