

Data Path : Z:\HPCHEM1\BNA E\DATA\BE041116\
 Data File : BE091633.D
 Acq On : 11 Apr 2016 18:02
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 12 15:35:55 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE040816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 08 16:48:14 2016
 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	120	0.00
2	1,4-Dioxane	0.640	0.618	3.4	114	0.00
3	n-Nitrosodimethylamine	0.767	0.688	10.3	116	0.00
4 S	2-Fluorophenol	1.193	1.018	14.7	108	0.00
5 S	Phenol-d6	1.591	1.376	13.5	114	0.00
6	bis(2-Chloroethyl)ether	1.379	1.310	5.0	120	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	122	-0.02
8 S	Nitrobenzene-d5	0.290	0.239	17.6	114	0.00
9	Nitrobenzene	0.305	0.244	20.0	113	0.00
10	Naphthalene	1.031	1.022	0.9	123	0.00
11	Hexachlorobutadiene	0.152	0.145	4.6	118	0.00
12	2-Methylnaphthalene	0.658	0.638	3.0	123	-0.01
13 I	Acenaphthene-d10	1.000	1.000	0.0	123	0.00
14 S	2,4,6-Tribromophenol	0.138	0.099	28.3#	111	0.00
15 S	2-Fluorobiphenyl	1.390	1.361	2.1	123	0.00
16	Acenaphthylene	1.941	1.744	10.1	137	0.00
17	Acenaphthene	1.227	1.255	-2.3	131	-0.01
18	Fluorene	1.533	1.464	4.5	122	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	122	0.00
20	4-Bromophenyl-phenylether	0.176	0.165	6.2	118	0.00
21	Hexachlorobenzene	0.183	0.182	0.5	122	0.00
22	Pentachlorophenol	0.068	0.049	27.9#	126	0.00
23	Phenanthrene	1.141	1.163	-1.9	125	0.00
24	Anthracene	1.011	0.959	5.1	123	0.00
25	Fluoranthene	1.208	1.109	8.2	121	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	114	0.00
27	Benzydine	0.329	0.244	25.8#	124	-0.02
28	Pvrene	1.000	0.961	3.9	119	-0.02
29 S	Terphenyl-d14	0.624	0.674	-8.0	126	0.00
30	Benzo(a)anthracene	1.119	1.096	2.1	116	0.00
31	3,3'-Dichlorobenzidine	0.599	0.380	36.6#	87	0.00
32	Chrysene	1.015	1.152	-13.5	132	-0.02
33	Bis(2-ethylhexyl)phthalate	0.494	0.178	64.0#	63	0.00
34	Indeno(1,2,3-cd)pyrene	1.132	0.985	13.0	103	-0.01
35 I	Perylene-d12	1.000	1.000	0.0	111	-0.01
36	Benzo(b)fluoranthene	1.161	1.053	9.3	106	-0.01
37	Benzo(k)fluoranthene	1.145	1.157	-1.0	118	-0.01
38 C	Benzo(a)pyrene	1.079	0.927	14.1	102	0.00
39	Dibenzo(a,h)anthracene	1.062	0.937	11.8	103	-0.01
40	Benzo(a,h,i)perylene	1.094	0.986	9.9	104	-0.01

Data Path : Z:\HPCHEM1\BNA E\DATA\BE041116\
Data File : BE091633.D
Acq On : 11 Apr 2016 18:02
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Apr 12 15:35:55 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE040816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 08 16:48:14 2016
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				