

Data Path : Z:\HPCHEM1\BNA M\DATA\BM020216\
 Data File : BM004171.D
 Acq On : 02 Feb 2016 19:34
 Operator : SJ/IZ
 Sample : SSTDC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDC001

Quant Time: Feb 03 01:15:51 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM012216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 22 14:46:41 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	138	0.00
2	1,4-Dioxane	0.547	0.404	26.1#	120	0.00
3	n-Nitrosodimethylamine	0.446	0.409	8.3	138	-0.02
4 S	2-Fluorophenol	1.043	0.983	5.8	143	-0.02
5 S	Phenol-d6	1.259	1.271	-1.0	155#	0.00
6	bis(2-Chloroethyl)ether	1.167	1.089	6.7	142	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	146	0.00
8 S	Nitrobenzene-d5	0.219	0.223	-1.8	165#	0.00
9	Nitrobenzene	0.250	0.254	-1.6	167#	0.00
10	Naphthalene	1.165	1.045	10.3	143	0.00
11	Hexachlorobutadiene	0.180	0.162	10.0	144	0.00
12	2-Methylnaphthalene	0.762	0.693	9.1	146	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	148	0.00
14 S	2,4,6-Tribromophenol	0.145	0.143	1.4	168#	0.00
15 S	2-Fluorobiphenyl	1.532	1.338	12.7	141	0.00
16	Acenaphthylene	2.158	2.079	3.7	158#	0.00
17	Acenaphthene	1.634	1.344	17.7	137	0.00
18	Fluorene	1.987	1.655	16.7	138	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	129	0.00
20	4-Bromophenyl-phenylether	0.163	0.193	-18.4	157#	0.00
21	Hexachlorobenzene	0.171	0.195	-14.0	154#	0.00
22	Pentachlorophenol	0.043	0.059	-37.2#	240#	0.00
23	Phenanthrene	1.074	1.171	-9.0	153#	0.00
24	Anthracene	1.142	1.196	-4.7	145	0.00
25	Fluoranthene	1.355	1.374	-1.4	141	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	142	0.00
27	Benzydine	0.241	0.301	-24.9	202#	0.00
28	Pvrene	1.423	1.299	8.7	144	0.00
29 S	Terphenyl-d14	0.643	0.588	8.6	147	0.00
30	Benzo(a)anthracene	1.518	1.322	12.9	140	0.00
31	3,3'-Dichlorobenzidine	0.338	0.337	0.3	173#	0.00
32	Chrysene	1.372	1.210	11.8	143	0.00
33	Bis(2-ethylhexyl)phthalate	0.595	0.606	-1.8	206#	-0.02
34	Indeno(1,2,3-cd)pyrene	1.463	1.156	21.0	127	0.00
35 I	Perylene-d12	1.000	1.000	0.0	131	0.00
36	Benzo(b)fluoranthene	1.551	1.400	9.7	127	-0.01
37	Benzo(k)fluoranthene	1.404	1.383	1.5	146	-0.01
38 C	Benzo(a)pyrene	1.366	1.308	4.2	141	0.00
39	Dibenzo(a,h)anthracene	1.227	1.079	12.1	128	-0.01
40	Benzo(a,h,i)perylene	1.339	1.119	16.4	123	-0.01

Data Path : Z:\HPCHEM1\BNA M\DATA\BM020216\
Data File : BM004171.D
Acq On : 02 Feb 2016 19:34
Operator : SJ/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC001

Quant Time: Feb 03 01:15:51 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM012216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 22 14:46:41 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				