

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012716\
 Data File : BM004138.D
 Acq On : 27 Jan 2016 17:25
 Operator : SJ/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Jan 27 23:59:39 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)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	74	0.00
2	1,4-Dioxane	0.547	0.475	13.2	76	0.00
3	n-Nitrosodimethylamine	0.446	0.407	8.7	74	0.00
4 S	2-Fluorophenol	1.043	0.953	8.6	75	0.00
5 S	Phenol-d6	1.259	1.168	7.2	77	0.00
6	bis(2-Chloroethyl)ether	1.167	1.080	7.5	76	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	79	0.00
8 S	Nitrobenzene-d5	0.219	0.202	7.8	80	0.00
9	Nitrobenzene	0.250	0.225	10.0	80	0.00
10	Naphthalene	1.165	1.064	8.7	78	0.00
11	Hexachlorobutadiene	0.180	0.162	10.0	78	0.00
12	2-Methylnaphthalene	0.762	0.719	5.6	81	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	86	0.00
14 S	2,4,6-Tribromophenol	0.145	0.132	9.0	90	0.00
15 S	2-Fluorobiphenyl	1.532	1.366	10.8	83	0.00
16	Acenaphthylene	2.158	1.913	11.4	84	0.00
17	Acenaphthene	1.634	1.433	12.3	85	0.00
18	Fluorene	1.987	1.852	6.8	90	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	92	0.00
20	4-Bromophenyl-phenylether	0.163	0.145	11.0	84	0.00
21	Hexachlorobenzene	0.171	0.152	11.1	87	0.00
22	Pentachlorophenol	0.043	0.036	16.3	105	0.00
23	Phenanthrene	1.074	0.990	7.8	92	0.00
24	Anthracene	1.142	1.038	9.1	90	0.00
25	Fluoranthene	1.355	1.274	6.0	94	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	86	0.00
27	Benzydine	0.241	0.217	10.0	88	0.02
28	Pyrene	1.423	1.341	5.8	90	0.00
29 S	Terphenyl-d14	0.643	0.621	3.4	94	0.00
30	Benzo(a)anthracene	1.518	1.376	9.4	88	0.00
31	3,3'-Dichlorobenzidine	0.338	0.330	2.4	103	0.00
32	Chrysene	1.372	1.436	-4.7	103	0.00
33	Bis(2-ethylhexyl)phthalate	0.595	0.526	11.6	109	0.00
34	Indeno(1,2,3-cd)pyrene	1.463	1.409	3.7	94	0.00
35 I	Perylene-d12	1.000	1.000	0.0	94	0.00
36	Benzo(b)fluoranthene	1.551	1.390	10.4	90	0.00
37	Benzo(k)fluoranthene	1.404	1.316	6.3	99	0.00
38 C	Benzo(a)pyrene	1.366	1.242	9.1	96	0.01
39	Dibenzo(a,h)anthracene	1.227	1.108	9.7	94	0.00
40	Benzo(g,h,i)perylene	1.339	1.206	9.9	94	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012716\
Data File : BM004138.D
Acq On : 27 Jan 2016 17:25
Operator : SJ/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC001

Quant Time: Jan 27 23:59:39 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				