

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030116\
 Data File : BM004510.D
 Acq On : 02 Mar 2016 08:57
 Operator : SJ/UM
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM030116

Quant Time: Mar 02 15:08:33 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM030116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 02 14:46:57 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	110	0.00
2	1,4-Dioxane	0.549	0.449	18.2	110	0.00
3	n-Nitrosodimethylamine	0.351	0.328	6.6	113	0.00
4 S	2-Fluorophenol	1.045	0.949	9.2	113	-0.02
5 S	Phenol-d6	1.208	1.118	7.5	111	-0.02
6	bis(2-Chloroethyl)ether	1.130	1.007	10.9	112	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
8 S	Nitrobenzene-d5	0.227	0.211	7.0	108	0.00
9	Nitrobenzene	0.249	0.234	6.0	110	0.00
10	Naphthalene	1.183	1.091	7.8	109	0.00
11	Hexachlorobutadiene	0.204	0.189	7.4	108	0.00
12	2-Methylnaphthalene	0.713	0.659	7.6	109	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
14 S	2,4,6-Tribromophenol	0.152	0.122	19.7	107	0.00
15 S	2-Fluorobiphenyl	1.573	1.442	8.3	107	0.00
16	Acenaphthylene	2.321	2.090	10.0	110	0.00
17	Acenaphthene	1.496	1.311	12.4	108	0.00
18	Fluorene	1.781	1.653	7.2	108	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
20	4-Bromophenyl-phenylether	0.201	0.168	16.4	108	0.00
21	Hexachlorobenzene	0.219	0.206	5.9	111	0.00
22	Pentachlorophenol	0.031	0.021	32.3#	102	0.00
23	Phenanthrene	1.222	1.121	8.3	109	0.00
24	Anthracene	1.096	0.910	17.0	107	0.00
25	Fluoranthene	1.591	1.312	17.5	110	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	107	0.02
27	Benzydine	0.420	0.374	11.0	107	0.00
28	Pvrene	1.949	1.729	11.3	111	0.00
29 S	Terphenyl-d14	0.966	0.798	17.4	109	0.00
30	Benzo(a)anthracene	1.680	1.380	17.9	106	0.02
31	3,3'-Dichlorobenzidine	0.478	0.472	1.3	113	0.00
32	Chrysene	2.064	1.817	12.0	122	0.00
33	Bis(2-ethylhexyl)phthalate	1.020	0.733	28.1#	96	0.00
34	Indeno(1,2,3-cd)pyrene	1.465	1.465	0.0	115	0.00
35 I	Perylene-d12	1.000	1.000	0.0	112	0.00
36	Benzo(b)fluoranthene	1.910	1.367	28.4#	106	0.00
37	Benzo(k)fluoranthene	1.727	1.524	11.8	120	0.00
38 C	Benzo(a)pyrene	1.484	1.348	9.2	113	0.00
39	Dibenzo(a,h)anthracene	1.227	1.112	9.4	113	0.00
40	Benzo(a,h,i)perylene	1.393	1.268	9.0	113	0.00

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030116\
Data File : BM004510.D
Acq On : 02 Mar 2016 08:57
Operator : SJ/UM
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM030116

Quant Time: Mar 02 15:08:33 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM030116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 02 14:46:57 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				