

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091015\
 Data File : BE090781.D
 Acq On : 11 Sep 2015 3:13
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Sep 11 03:55:46 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 11:40:50 2015
 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	110	0.00
2	1,4-Dioxane	0.814	0.611	24.9	85	0.00
3	n-Nitrosodimethylamine	1.033	0.752	27.2#	85	0.00
4 S	2-Fluorophenol	1.039	0.830	20.1	94	0.00
5 S	Phenol-d6	1.392	1.249	10.3	104	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	121	0.00
7 S	Nitrobenzene-d5	0.330	0.274	17.0	101	0.00
8	Nitrobenzene	0.365	0.306	16.2	106	0.00
9	Naphthalene	1.091	0.938	14.0	105	0.00
10	Hexachlorobutadiene	0.171	0.144	15.8	99	0.00
11	2-Methylnaphthalene	0.590	0.534	9.5	113	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	132	0.00
13 S	2,4,6-Tribromophenol	0.143	0.106	25.9#	116	0.00
14 S	2-Fluorobiphenyl	1.385	1.139	17.8	111	0.00
15	Acenaphthylene	7.992	6.789	15.1	117	0.00
16	Acenaphthene	1.335	1.131	15.3	111	0.00
17	Fluorene	1.666	1.423	14.6	114	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	127	0.00
19	4-Bromophenyl-phenylether	0.170	0.143	15.9	111	0.00
20	Hexachlorobenzene	0.928	0.763	17.8	106	0.00
21	Pentachlorophenol	0.054	0.037	31.5#	99	0.00
22	Phenanthrene	1.237	1.057	14.6	109	-0.02
23	Anthracene	1.035	0.908	12.3	116	0.00
24	Fluoranthene	1.304	1.077	17.4	114	0.00
25 I	Chrysene-d12	1.000	1.000	0.0	121	0.00
26	Pyrene	1.035	0.949	8.3	117	0.00
27 S	Terphenyl-d14	0.553	0.493	10.8	113	0.00
28	Benzo(a)anthracene	1.156	0.986	14.7	109	0.00
29	Chrysene	1.279	1.113	13.0	107	0.00
30	Bis(2-ethylhexyl)phthalate	0.345	0.226	34.5#	108	0.00
31	Indeno(1,2,3-cd)pyrene	1.181	1.032	12.6	115	0.00
32 I	Perylene-d12	1.000	1.000	0.0	125	0.00
33	Benzo(b)fluoranthene	1.084	0.917	15.4	109	0.00
34	Benzo(k)fluoranthene	1.263	1.116	11.6	112	0.00
35 C	Benzo(a)pyrene	0.970	0.846	12.8	117	0.00
36	Dibenzo(a,h)anthracene	1.020	0.890	12.7	114	0.00
37	Benzo(g,h,i)perylene	1.116	0.924	17.2	110	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0