

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE082815\
 Data File : BE090656.D
 Acq On : 28 Aug 2015 11:51
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Aug 31 04:57:43 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE081415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 14 16:52:42 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	103	0.00
2	1,4-Dioxane	0.658	0.757	-15.0	112	0.00
3	n-Nitrosodimethylamine	0.863	0.852	1.3	103	0.00
4 S	2-Fluorophenol	1.446	0.894	38.2#	63	0.00
5 S	Phenol-d6	2.125	1.333	37.3#	67	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	94	0.00
7 S	Nitrobenzene-d5	0.398	0.353	11.3	88	0.00
8	Nitrobenzene	0.413	0.348	15.7	85	0.00
9	Naphthalene	1.136	1.124	1.1	94	0.00
10	Hexachlorobutadiene	0.167	0.181	-8.4	102	0.00
11	2-Methylnaphthalene	0.740	0.659	10.9	86	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
13 S	2,4,6-Tribromophenol	0.192	0.121	37.0#	71	0.00
14 S	2-Fluorobiphenyl	1.493	1.316	11.9	90	0.00
15	Acenaphthylene	9.132	8.055	11.8	92	0.00
16	Acenaphthene	1.374	1.349	1.8	102	0.00
17	Fluorene	1.684	1.762	-4.6	108	0.02
18 I	Phenanthrene-d10	1.000	1.000	0.0	120	0.00
19	4-Bromophenyl-phenylether	0.204	0.183	10.3	111	0.00
20	Hexachlorobenzene	0.906	0.912	-0.7	121	-0.02
21	Pentachlorophenol	0.079	0.040	49.4#	80	0.00
22	Phenanthrene	1.312	1.257	4.2	117	0.00
23	Anthracene	1.208	1.073	11.2	110	0.00
24	Fluoranthene	1.396	1.276	8.6	113	0.00
25 I	Chrysene-d12	1.000	1.000	0.0	129	0.00
26	Pyrene	1.430	1.253	12.4	116	0.00
27 S	Terphenyl-d14	0.914	0.788	13.8	113	0.00
28	Benzo(a)anthracene	1.316	1.174	10.8	121	0.00
29	Chrysene	1.344	1.334	0.7	130	0.00
30	Bis(2-ethylhexyl)phthalate	0.725	0.288	60.3#	62	0.00
31	Indeno(1,2,3-cd)pyrene	1.266	1.006	20.5	111	0.00
32 I	Perylene-d12	1.000	1.000	0.0	120	0.00
33	Benzo(b)fluoranthene	1.184	1.110	6.2	116	0.00
34	Benzo(k)fluoranthene	1.342	1.339	0.2	119	0.00
35 C	Benzo(a)pyrene	1.108	0.984	11.2	110	0.00
36	Dibenzo(a,h)anthracene	1.069	0.881	17.6	105	0.00
37	Benzo(g,h,i)perylene	1.169	1.001	14.4	105	0.00

(#) = Out of Range

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