

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070915\
 Data File : BE090079.D
 Acq On : 9 Jul 2015 18:46
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Jul 10 02:19:48 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE062915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 30 07:01:01 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	125	-0.01
2	1,4-Dioxane	0.686	0.546	20.4	103	0.00
3	n-Nitrosodimethylamine	0.930	0.724	22.2	101	-0.01
4 S	2-Fluorophenol	1.149	1.188	-3.4	129	0.00
5 S	Phenol-d6	1.608	1.949	-21.2	154#	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	134	-0.01
7 S	Nitrobenzene-d5	0.397	0.357	10.1	123	-0.01
8	Nitrobenzene	0.421	0.343	18.5	113	0.00
9	Naphthalene	1.130	0.979	13.4	118	-0.01
10	Hexachlorobutadiene	0.178	0.158	11.2	121	0.00
11	2-Methylnaphthalene	0.755	0.628	16.8	114	-0.01
12 I	Acenaphthene-d10	1.000	1.000	0.0	125	-0.01
13 S	2,4,6-Tribromophenol	0.147	0.157	-6.8	137	0.00
14 S	2-Fluorobiphenyl	1.503	1.448	3.7	124	-0.01
15	Acenaphthylene	9.019	7.717	14.4	110	0.00
16	Acenaphthene	1.300	1.132	12.9	110	-0.01
17	Fluorene	1.768	1.416	19.9	103	-0.02
18 I	Phenanthrene-d10	1.000	1.000	0.0	109	-0.02
19	4-Bromophenyl-phenylether	0.213	0.179	16.0	94	-0.02
20	Hexachlorobenzene	0.885	0.772	12.8	99	-0.02
21	Pentachlorophenol	0.086	0.072	16.3	96	0.00
22	Phenanthrene	1.281	1.257	1.9	111	-0.02
23	Anthracene	1.172	1.012	13.7	96	0.00
24	Fluoranthene	1.431	1.564	-9.3	122	-0.02
25 I	Chrysene-d12	1.000	1.000	0.0	97	-0.01
26	Pyrene	1.192	1.404	-17.8	116	-0.01
27 S	Terphenyl-d14	0.833	0.838	-0.6	97	-0.01
28	Benzo(a)anthracene	1.249	1.266	-1.4	100	-0.01
29	Chrysene	1.297	1.241	4.3	94	-0.01
30	Bis(2-ethylhexyl)phthalate	0.398	0.453	-13.8	120	-0.01
31	Indeno(1,2,3-cd)pyrene	1.243	1.167	6.1	92	-0.04
32 I	Perylene-d12	1.000	1.000	0.0	92	-0.02
33	Benzo(b)fluoranthene	1.103	1.220	-10.6	102	-0.02
34	Benzo(k)fluoranthene	1.232	1.257	-2.0	94	-0.02
35 C	Benzo(a)pyrene	1.019	1.117	-9.6	101	-0.02
36	Dibenzo(a,h)anthracene	1.046	0.972	7.1	87	-0.03
37	Benzo(g,h,i)perylene	1.060	1.054	0.6	92	-0.04

(#) = Out of Range

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