

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080117\
 Data File : BE093618.D
 Acq On : 1 Aug 2017 16:04
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE080117

Quant Time: Aug 01 17:33:28 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 01 16:48:54 2017
 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	113	0.00
2	1,4-Dioxane	0.648	0.630	2.8	110	0.01
3	n-Nitrosodimethylamine	0.607	0.579	4.6	114	0.01
4 S	2-Fluorophenol	1.192	1.155	3.1	114	0.00
5 S	Phenol-d6	1.794	1.686	6.0	110	0.00
6	bis(2-Chloroethyl)ether	1.508	1.465	2.9	112	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	109	0.00
8 S	Nitrobenzene-d5	0.305	0.293	3.9	112	0.00
9	Naphthalene	4.630	4.510	2.6	107	0.00
10	Hexachlorobutadiene	0.172	0.167	2.9	108	0.00
11 SURR	2-Methylnaphthalene-d10	0.632	0.618	2.2	107	0.00
12	2-Methylnaphthalene	0.766	0.736	3.9	105	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
14 S	2,4,6-Tribromophenol	0.093	0.072	22.6	102	0.00
15 S	2-Fluorobiphenyl	1.601	1.571	1.9	105	-0.01
16	Acenaphthylene	1.912	1.807	5.5	105	0.00
17	Acenaphthene	1.349	1.290	4.4	104	0.00
18	Fluorene	1.803	1.705	5.4	100	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	113	0.00
20	4-Bromophenyl-phenylether	0.126	0.120	4.8	97	0.00
21	Hexachlorobenzene	0.129	0.133	-3.1	108	0.00
22	Pentachlorophenol	0.055	0.048	12.7	103	0.00
23	Phenanthrene	0.880	0.845	4.0	106	0.00
24	Anthracene	1.203	1.151	4.3	102	0.00
25 SURR	Fluoranthene-d10	4.541	4.410	2.9	101	0.00
26	Fluoranthene	1.401	1.364	2.6	101	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	103	0.00
28	Pyrene	1.292	1.239	4.1	99	0.00
29 S	Terphenyl-d14	0.719	0.709	1.4	101	0.00
30	Benzo(a)anthracene	1.176	1.134	3.6	103	0.00
31	Chrysene	1.311	1.286	1.9	101	0.00
32	Bis(2-ethylhexyl)phthalate	1.603	1.354	15.5	101	0.00
33	Indeno(1,2,3-cd)pyrene	1.187	1.161	2.2	100	0.00
34 I	Perylene-d12	1.000	1.000	0.0	102	0.00
35	Benzo(b)fluoranthene	1.427	1.364	4.4	100	0.00
36	Benzo(k)fluoranthene	1.459	1.427	2.2	103	0.00
37 C	Benzo(a)pyrene	1.293	1.238	4.3	101	0.00
38	Dibenzo(a,h)anthracene	1.248	1.224	1.9	101	0.00
39	Benzo(g,h,i)perylene	1.266	1.234	2.5	101	0.00

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(#) = Out of Range	SPCC's out = 0	CCC's out = 0		