

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100617\
 Data File : BE094180.D
 Acq On : 6 Oct 2017 15:07
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICSVBE100617

Quant Time: Oct 06 15:47:25 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE100617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 06 15:16:13 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	104	0.00
2	1,4-Dioxane	0.725	0.685	5.5	92	0.02
3	n-Nitrosodimethylamine	0.777	0.744	4.2	100	0.02
4 S	2-Fluorophenol	1.333	1.306	2.0	102	0.00
5 S	Phenol-d6	1.932	1.884	2.5	103	0.00
6	bis(2-Chloroethyl)ether	1.569	1.557	0.8	105	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
8 S	Nitrobenzene-d5	0.333	0.330	0.9	107	0.00
9	Naphthalene	4.505	4.558	-1.2	106	0.00
10	Hexachlorobutadiene	0.171	0.174	-1.8	108	-0.02
11 SURR	2-Methylnaphthalene-d10	0.616	0.624	-1.3	107	0.00
12	2-Methylnaphthalene	0.740	0.748	-1.1	107	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
14 S	2,4,6-Tribromophenol	0.174	0.142	18.4	95	0.00
15 S	2-Fluorobiphenyl	1.370	1.360	0.7	108	0.00
16	Acenaphthylene	2.026	1.980	2.3	106	0.00
17	Acenaphthene	1.338	1.333	0.4	106	0.00
18	Fluorene	1.753	1.727	1.5	107	-0.02
19 I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
20	4-Bromophenyl-phenylether	0.230	0.224	2.6	101	0.00
21	Hexachlorobenzene	0.207	0.193	6.8	99	0.00
22	Pentachlorophenol	0.056	0.048	14.3	89	0.00
23	Phenanthrene	1.386	1.348	2.7	101	0.00
24	Anthracene	1.057	1.082	-2.4	110	0.00
25 SURR	Fluoranthene-d10	3.995	3.746	6.2	97	0.00
26	Fluoranthene	1.216	1.143	6.0	98	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	95	0.00
28	Pvrene	1.376	1.371	0.4	95	0.00
29 S	Terphenyl-d14	0.755	0.752	0.4	95	0.00
30	Benzo(a)anthracene	1.314	1.287	2.1	93	0.00
31	Chrysene	1.297	1.315	-1.4	97	0.00
32	Bis(2-ethylhexyl)phthalate	3.027	2.687	11.2	89	0.00
33	Indeno(1,2,3-cd)pyrene	1.287	1.308	-1.6	96	0.00
34 I	Perylene-d12	1.000	1.000	0.0	95	0.00
35	Benzo(b)fluoranthene	1.345	1.358	-1.0	96	0.00
36	Benzo(k)fluoranthene	1.368	1.380	-0.9	96	0.00
37 C	Benzo(a)pyrene	1.272	1.262	0.8	95	0.00
38	Dibenzo(a,h)anthracene	1.179	1.171	0.7	94	0.00
39	Benzo(a,h,i)perylene	1.204	1.216	-1.0	97	-0.01

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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