

Data Path : Z:\HPCHEM1\BNA E\DATA\BE092317\
 Data File : BE094079.D
 Acq On : 23 Sep 2017 15:45
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE092317

Quant Time: Sep 25 13:07:06 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 25 13:07:29 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	92	0.00
2	1,4-Dioxane	0.835	1.131	-35.4#	126	0.00
3	n-Nitrosodimethylamine	1.034	0.917	11.3	98	0.00
4 S	2-Fluorophenol	2.100	2.151	-2.4	94	0.00
5 S	Phenol-d6	2.365	2.309	2.4	90	0.00
6	bis(2-Chloroethyl)ether	1.767	1.699	3.8	88	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
8 S	Nitrobenzene-d5	0.356	0.343	3.7	89	0.00
9	Naphthalene	5.096	4.656	8.6	87	0.00
10	Hexachlorobutadiene	0.162	0.163	-0.6	93	0.00
11 SURR	2-Methylnaphthalene-d10	0.678	0.650	4.1	90	0.00
12	2-Methylnaphthalene	0.808	0.777	3.8	91	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	89	0.00
14 S	2,4,6-Tribromophenol	0.303	0.133	56.1#	48#	0.00
15 S	2-Fluorobiphenyl	1.314	1.341	-2.1	90	0.00
16	Acenaphthylene	2.014	1.953	3.0	87	0.00
17	Acenaphthene	1.336	1.326	0.7	88	0.00
18	Fluorene	1.989	1.926	3.2	88	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	92	0.00
20	4-Bromophenyl-phenylether	0.214	0.197	7.9	90	0.00
21	Hexachlorobenzene	0.157	0.144	8.3	92	0.00
22	Pentachlorophenol	0.011	0.022	-100.0#	275#	0.00
23	Phenanthrene	1.433	1.358	5.2	91	0.00
24	Anthracene	1.105	1.083	2.0	87	0.00
25 SURR	Fluoranthene-d10	5.132	4.874	5.0	94	0.00
26	Fluoranthene	1.537	1.458	5.1	94	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	99	-0.01
28	Pvrene	1.107	1.123	-1.4	94	0.00
29 S	Terphenyl-d14	0.627	0.638	-1.8	95	0.00
30	Benzo(a)anthracene	1.340	1.343	-0.2	99	0.00
31	Chrysene	1.313	1.349	-2.7	100	0.00
32	Bis(2-ethylhexyl)phthalate	3.600	3.377	6.2	92	0.00
33	Indeno(1,2,3-cd)pyrene	1.015	0.971	4.3	97	0.00
34 I	Perylene-d12	1.000	1.000	0.0	98	0.00
35	Benzo(b)fluoranthene	1.448	1.501	-3.7	101	0.00
36	Benzo(k)fluoranthene	1.453	1.446	0.5	97	0.00
37 C	Benzo(a)pyrene	1.300	1.302	-0.2	99	0.00
38	Dibenzo(a,h)anthracene	1.041	1.060	-1.8	98	0.00
39	Benzo(a,h,i)perylene	1.040	1.065	-2.4	98	0.00

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