

Data Path : Z:\HPCHEM1\BNA E\Data\BE092517\
 Data File : BE094106.D
 Acq On : 25 Sep 2017 17:22
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE092517

Quant Time: Sep 25 17:58:08 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 25 17:37:08 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	98	0.00
2	1,4-Dioxane	0.780	0.653	16.3	79	0.00
3	n-Nitrosodimethylamine	0.996	0.854	14.3	84	0.00
4 S	2-Fluorophenol	1.516	1.431	5.6	91	0.00
5 S	Phenol-d6	2.106	1.962	6.8	90	0.00
6	bis(2-Chloroethyl)ether	1.632	1.558	4.5	92	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	91	0.00
8 S	Nitrobenzene-d5	0.387	0.367	5.2	92	0.00
9	Naphthalene	4.516	4.545	-0.6	93	0.00
10	Hexachlorobutadiene	0.169	0.168	0.6	95	0.00
11 SURR	2-Methylnaphthalene-d10	0.616	0.615	0.2	93	0.00
12	2-Methylnaphthalene	0.731	0.723	1.1	93	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	91	0.00
14 S	2,4,6-Tribromophenol	0.256	0.176	31.3#	74	0.00
15 S	2-Fluorobiphenyl	1.322	1.406	-6.4	93	0.02
16	Acenaphthylene	2.005	2.012	-0.3	89	0.00
17	Acenaphthene	1.253	1.331	-6.2	91	0.00
18	Fluorene	1.700	1.737	-2.2	91	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	87	0.00
20	4-Bromophenyl-phenylether	0.164	0.163	0.6	89	0.00
21	Hexachlorobenzene	0.104	0.114	-9.6	97	0.00
22	Pentachlorophenol	0.095	0.086	9.5	82	0.00
23	Phenanthrene	1.023	1.031	-0.8	91	0.00
24	Anthracene	1.074	1.070	0.4	92	0.00
25 SURR	Fluoranthene-d10	3.873	3.962	-2.3	93	0.00
26	Fluoranthene	1.147	1.189	-3.7	93	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	98	0.00
28	Pvrene	1.267	1.174	7.3	94	0.00
29 S	Terphenyl-d14	0.730	0.666	8.8	94	0.00
30	Benzo(a)anthracene	1.347	1.338	0.7	98	0.00
31	Chrysene	1.300	1.311	-0.8	99	0.00
32	Bis(2-ethylhexyl)phthalate	6.447	4.269	33.8#	87	0.00
33	Indeno(1,2,3-cd)pyrene	1.391	1.375	1.2	101	0.00
34 I	Perylene-d12	1.000	1.000	0.0	99	0.00
35	Benzo(b)fluoranthene	1.443	1.342	7.0	102	0.00
36	Benzo(k)fluoranthene	1.382	1.334	3.5	95	0.00
37 C	Benzo(a)pyrene	1.304	1.270	2.6	100	0.00
38	Dibenzo(a,h)anthracene	1.235	1.208	2.2	102	0.00
39	Benzo(a,h,i)perylene	1.222	1.208	1.1	102	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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