

Data Path : Z:\HPCHEM1\BNA E\DATA\BE041816\
 Data File : BE091667.D
 Acq On : 18 Apr 2016 16:19
 Operator : UM/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE041816

Quant Time: Apr 18 17:40:25 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE041816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 18 17:20:40 2016
 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	92	0.00
2	1,4-Dioxane	0.400	0.401	-0.3	99	0.00
3	n-Nitrosodimethylamine	0.400	0.428	-7.0	106	0.00
4 S	2-Fluorophenol	0.400	0.408	-2.0	99	0.00
5 S	Phenol-d6	0.400	0.388	3.0	97	0.00
6	bis(2-Chloroethyl)ether	0.400	0.408	-2.0	98	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	91	0.00
8 S	Nitrobenzene-d5	0.400	0.404	-1.0	107	0.00
9	Nitrobenzene	0.400	0.406	-1.5	105	0.00
10	Naphthalene	0.400	0.413	-3.2	94	0.00
11	Hexachlorobutadiene	0.400	0.428	-7.0	99	0.00
12	2-Methylnaphthalene	0.400	0.404	-1.0	95	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	92	0.00
14 S	2,4,6-Tribromophenol	0.400	0.449	-12.2	104	-0.01
15 S	2-Fluorobiphenyl	0.400	0.411	-2.7	95	0.00
16	Acenaphthylene	0.400	0.404	-1.0	106	0.00
17	Acenaphthene	0.400	0.419	-4.7	93	0.00
18	Fluorene	0.400	0.401	-0.3	94	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	84	0.00
20	4-Bromophenyl-phenylether	0.400	0.401	-0.3	98	0.00
21	Hexachlorobenzene	0.400	0.398	0.5	92	0.00
22	Pentachlorophenol	0.400	0.451	-12.7	101	0.00
23	Phenanthrene	0.400	0.414	-3.5	95	0.00
24	Anthracene	0.400	0.390	2.5	96	0.00
25	Fluoranthene	0.400	0.413	-3.2	101	0.00
26 I	Chrysene-d12	5.000	5.000	0.0	97	0.00
27	Benzydine	0.400	0.412	-3.0	124	-0.02
28	Pvrene	0.400	0.465	-16.3	120	0.00
29 S	Terphenyl-d14	0.400	0.412	-3.0	105	0.00
30	Benzo(a)anthracene	0.400	0.432	-8.0	110	0.00
31	3,3'-Dichlorobenzidine	0.400	0.464	-16.0	125	0.00
32	Chrysene	0.400	0.405	-1.3	97	0.00
33	Bis(2-ethylhexyl)phthalate	0.400	0.547	-36.8#	141	0.00
34	Indeno(1,2,3-cd)pyrene	0.400	0.421	-5.2	106	0.00
35 I	Perylene-d12	5.000	5.000	0.0	95	-0.01
36	Benzo(b)fluoranthene	0.400	0.423	-5.7	108	0.00
37	Benzo(k)fluoranthene	0.400	0.420	-5.0	105	-0.01
38 C	Benzo(a)pyrene	0.400	0.422	-5.5	109	0.00
39	Dibenzo(a,h)anthracene	0.400	0.421	-5.2	107	0.00
40	Benzo(a,h,i)perylene	0.400	0.408	-2.0	102	0.00

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Compound	Amount	Calc.	%Dev Area	% Dev(min)

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