

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072716\
 Data File : BE092058.D
 Acq On : 27 Jul 2016 15:52
 Operator : UM/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE072716

Quant Time: Jul 27 17:39:13 2016
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE072716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 27 17:13:26 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	-0.02
2	1,4-Dioxane	0.400	0.410	-2.5	95	0.00
3	n-Nitrosodimethylamine	0.400	0.366	8.5	93	0.00
4 S	2-Fluorophenol	0.400	0.356	11.0	94	-0.02
5 S	Phenol-d6	0.400	0.325	18.8	85	0.00
6	bis(2-Chloroethyl)ether	0.400	0.359	10.3	92	0.00
7	n-Nitroso-di-n-propylamine	0.400	0.356	11.0	93	0.00
8 I	Naphthalene-d8	5.000	5.000	0.0	92	0.00
9 S	Nitrobenzene-d5	0.400	0.328	18.0	81	0.00
10	Nitrobenzene	0.400	0.351	12.3	91	0.00
11	bis(2-Chloroethoxy)methane	0.400	0.347	13.3	92	0.00
12	Naphthalene	0.400	0.392	2.0	91	0.00
13	Hexachlorobutadiene	0.400	0.399	0.3	94	0.00
14	2-Methylnaphthalene	0.400	0.377	5.8	92	0.00
15 I	Acenaphthene-d10	5.000	5.000	0.0	87	0.00
16 S	2,4,6-Tribromophenol	0.400	0.431	-7.7	89	0.00
17 S	2-Fluorobiphenyl	0.400	0.409	-2.2	91	0.00
18	Acenaphthylene	0.400	0.387	3.3	92	0.00
19	2,6-Dinitrotoluene	0.400	0.423	-5.7	87	0.00
20	Acenaphthene	0.400	0.399	0.3	89	0.00
21	2,4-Dinitrotoluene	0.400	0.336	16.0	81	0.00
22	Fluorene	0.400	0.393	1.8	89	0.00
23	4-Chlorophenyl-phenylether	0.400	0.419	-4.7	92	0.00
24 I	Phenanthrene-d10	5.000	5.000	0.0	89	0.00
25	4-Bromophenyl-phenylether	0.400	0.384	4.0	88	0.00
26	Hexachlorobenzene	0.400	0.394	1.5	89	0.00
27	Pentachlorophenol	0.400	0.467	-16.8	82	0.00
28	Phenanthrene	0.400	0.396	1.0	91	0.00
29	Anthracene	0.400	0.370	7.5	88	0.00
30	Fluoranthene	0.400	0.365	8.8	92	0.00
31 I	Chrysene-d12	5.000	5.000	0.0	79	0.00
32	Benzidine	0.400	0.408	-2.0	85	0.00
33	Pyrene	0.400	0.415	-3.7	90	0.00
34 S	Terphenyl-d14	0.400	0.401	-0.3	89	0.00
35	Benzo(a)anthracene	0.400	0.438	-9.5	86	0.00
36	3,3'-Dichlorobenzidine	0.400	0.463	-15.8	90	0.00
37	Chrysene	0.400	0.328	18.0	79	0.00
38	Bis(2-ethylhexyl)phthalate	0.400	0.480	-20.0	98	0.00
39	Indeno(1,2,3-cd)pyrene	0.400	0.450	-12.5	96	0.00
40 I	Perylene-d12	5.000	5.000	0.0	93	0.00
41	Benzo(b)fluoranthene	0.400	0.410	-2.5	97	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.413	-3.2	91	0.00
43 C	Benzo(a)pyrene	0.400	0.377	5.8	90	0.00
44	Dibenzo(a,h)anthracene	0.400	0.382	4.5	96	0.00
45	Benzo(g,h,i)perylene	0.400	0.390	2.5	97	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0