

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071615\
 Data File : BE090133.D
 Acq On : 16 Jul 2015 23:54
 Operator : TP/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE071615

Quant Time: Jul 17 13:00:54 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE071615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 17 12:59:41 2015
 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	87	0.00
2	1,4-Dioxane	1.000	1.060	-6.0	90	0.00
3	n-Nitrosodimethylamine	1.000	0.950	5.0	87	0.00
4 S	2-Fluorophenol	1.000	0.996	0.4	93	0.00
5 S	Phenol-d6	1.000	1.021	-2.1	93	0.00
6 I	Naphthalene-d8	5.000	5.000	0.0	88	0.00
7 S	Nitrobenzene-d5	1.000	0.983	1.7	90	0.00
8	Nitrobenzene	1.000	0.993	0.7	91	0.00
9	Naphthalene	1.000	0.971	2.9	89	0.00
10	Hexachlorobutadiene	1.000	0.958	4.2	88	0.00
11	2-Methylnaphthalene	1.000	0.994	0.6	91	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	90	0.00
13 S	2,4,6-Tribromophenol	1.000	0.919	8.1	106	0.00
14 S	2-Fluorobiphenyl	1.000	0.971	2.9	89	0.00
15	Acenaphthylene	1.000	0.980	2.0	91	0.00
16	Acenaphthene	1.000	0.967	3.3	90	0.00
17	Fluorene	1.000	0.963	3.7	89	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	92	0.00
19	4-Bromophenyl-phenylether	1.000	0.975	2.5	93	0.00
20	Hexachlorobenzene	1.000	0.963	3.7	89	0.00
21	Pentachlorophenol	1.000	1.038	-3.8	102	0.00
22	Phenanthrene	1.000	0.955	4.5	91	0.00
23	Anthracene	1.000	0.965	3.5	91	0.00
24	Fluoranthene	1.000	0.996	0.4	93	0.00
25 I	Chrysene-d12	5.000	5.000	0.0	97	0.00
26	Pyrene	1.000	0.928	7.2	92	0.00
27 S	Terphenyl-d14	1.000	0.989	1.1	95	0.00
28	Benzo(a)anthracene	1.000	1.070	-7.0	110	0.00
29	Chrysene	1.000	1.036	-3.6	97	0.00
30	Bis(2-ethylhexyl)phthalate	1.000	0.979	2.1	101	0.00
31	Indeno(1,2,3-cd)pyrene	1.000	1.178	-17.8	123	0.00
32 I	Perylene-d12	5.000	5.000	0.0	98	0.00
33	Benzo(b)fluoranthene	1.000	1.324	-32.4#	132	0.00
34	Benzo(k)fluoranthene	1.000	0.947	5.3	104	0.00
35 C	Benzo(a)pyrene	1.000	0.942	5.8	124	0.00
36	Dibenzo(a,h)anthracene	1.000	1.247	-24.7	133	0.00
37	Benzo(g,h,i)perylene	1.000	0.941	5.9	117	0.00

(#) = Out of Range

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