

Data Path : Z:\HPCHEM1\BNA E\DATA\BE051117\
 Data File : BE092995.D
 Acq On : 11 May 2017 14:27
 Operator : SJ/MA
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE051117

Quant Time: May 11 16:40:41 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE051117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 11 14:48:08 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	88	-0.02
2	1,4-Dioxane	0.400	0.389	2.8	101	0.00
3	n-Nitrosodimethylamine	0.400	0.393	1.8	100	0.00
4 S	2-Fluorophenol	0.400	0.387	3.3	99	0.00
5 S	Phenol-d6	0.400	0.395	1.3	101	-0.02
6	bis(2-Chloroethyl)ether	0.400	0.402	-0.5	102	-0.02
7 I	Naphthalene-d8	0.400	0.400	0.0	89	0.00
8 S	Nitrobenzene-d5	0.400	0.389	2.8	100	-0.02
9	Naphthalene	0.400	0.405	-1.3	100	0.00
10	Hexachlorobutadiene	0.400	0.407	-1.7	102	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.403	-0.8	102	-0.02
12	2-Methylnaphthalene	0.400	0.395	1.3	100	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	90	0.00
14 S	2,4,6-Tribromophenol	0.400	0.375	6.3	104	0.00
15 S	2-Fluorobiphenyl	0.400	0.395	1.3	101	0.00
16	Acenaphthylene	0.400	0.376	6.0	102	0.00
17	Acenaphthene	0.400	0.394	1.5	101	0.00
18	Fluorene	0.400	0.391	2.3	100	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	93	0.00
20	4-Bromophenyl-phenylether	0.400	0.383	4.3	101	0.00
21	Hexachlorobenzene	0.400	0.424	-6.0	105	0.00
22	Pentachlorophenol	0.400	0.433	-8.2	101	0.00
23	Phenanthrene	0.400	0.407	-1.7	104	0.01
24	Anthracene	0.400	0.423	-5.7	107	0.00
25 SURR	Fluoranthene-d10	0.400	0.393	1.8	102	-0.02
26	Fluoranthene	0.400	0.401	-0.3	104	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	94	0.00
28	Pyrene	0.400	0.412	-3.0	107	0.00
29 S	Terphenyl-d14	0.400	0.383	4.3	108	0.00
30	Benzo(a)anthracene	0.400	0.371	7.3	108	0.00
31	Chrysene	0.400	0.416	-4.0	109	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.390	2.5	107	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.428	-7.0	114	0.00
34 I	Perylene-d12	0.400	0.400	0.0	95	0.00
35	Benzo(b)fluoranthene	0.400	0.373	6.8	96	0.00
36	Benzo(k)fluoranthene	0.400	0.421	-5.2	125	0.00
37 C	Benzo(a)pyrene	0.400	0.378	5.5	106	0.00
38	Dibenzo(a,h)anthracene	0.400	0.426	-6.5	119	0.02
39	Benzo(a,h,i)perylene	0.400	0.416	-4.0	109	0.00

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

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