

Data Path : Z:\HPCHEM1\BNA E\DATA\BE092317\
 Data File : BE094079.D
 Acq On : 23 Sep 2017 15:45
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE092317

Quant Time: Sep 25 13:07:06 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 25 13:07:29 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	92	0.00
2	1,4-Dioxane	0.400	0.609	-52.2#	126	0.00
3	n-Nitrosodimethylamine	0.400	0.372	7.0	98	0.00
4 S	2-Fluorophenol	0.400	0.395	1.3	94	0.00
5 S	Phenol-d6	0.400	0.391	2.3	90	0.00
6	bis(2-Chloroethyl)ether	0.400	0.385	3.8	88	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	92	0.00
8 S	Nitrobenzene-d5	0.400	0.385	3.8	89	0.00
9	Naphthalene	0.400	0.365	8.8	87	0.00
10	Hexachlorobutadiene	0.400	0.402	-0.5	93	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.384	4.0	90	0.00
12	2-Methylnaphthalene	0.400	0.385	3.8	91	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	89	0.00
14 S	2,4,6-Tribromophenol	0.400	0.176	56.0#	48	0.00
15 S	2-Fluorobiphenyl	0.400	0.408	-2.0	90	0.00
16	Acenaphthylene	0.400	0.388	3.0	87	0.00
17	Acenaphthene	0.400	0.397	0.8	88	0.00
18	Fluorene	0.400	0.387	3.3	88	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	92	0.00
20	4-Bromophenyl-phenylether	0.400	0.369	7.8	90	0.00
21	Hexachlorobenzene	0.400	0.367	8.3	92	0.00
22	Pentachlorophenol	0.400	0.787	-96.8#	275	0.00
23	Phenanthrene	0.400	0.379	5.3	91	0.00
24	Anthracene	0.400	0.392	2.0	87	0.00
25 SURR	Fluoranthene-d10	0.400	0.380	5.0	94	0.00
26	Fluoranthene	0.400	0.379	5.3	94	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	99	-0.01
28	Pvrene	0.400	0.406	-1.5	94	0.00
29 S	Terphenyl-d14	0.400	0.407	-1.7	95	0.00
30	Benzo(a)anthracene	0.400	0.401	-0.3	99	0.00
31	Chrysene	0.400	0.411	-2.7	100	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.375	6.3	92	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.383	4.3	97	0.00
34 I	Perylene-d12	0.400	0.400	0.0	98	0.00
35	Benzo(b)fluoranthene	0.400	0.415	-3.7	101	0.00
36	Benzo(k)fluoranthene	0.400	0.398	0.5	97	0.00
37 C	Benzo(a)pyrene	0.400	0.401	-0.3	99	0.00
38	Dibenzo(a,h)anthracene	0.400	0.407	-1.7	98	0.00
39	Benzo(a,h,i)perylene	0.400	0.409	-2.2	98	0.00

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

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