

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102317\
 Data File : BE094501.D
 Acq On : 23 Oct 2017 18:33
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE102317

Quant Time: Oct 23 19:34:29 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 23 19:26:42 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	102	0.00
2	1,4-Dioxane	0.677	0.688	-1.6	89	0.00
3	n-Nitrosodimethylamine	0.648	0.601	7.3	94	0.00
4 S	2-Fluorophenol	1.370	1.403	-2.4	104	0.00
5 S	Phenol-d6	2.069	2.256	-9.0	114	0.00
6	bis(2-Chloroethyl)ether	1.600	1.591	0.6	102	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
8 S	Nitrobenzene-d5	0.319	0.314	1.6	106	0.00
9	Naphthalene	4.527	4.566	-0.9	101	0.00
10	Hexachlorobutadiene	0.164	0.167	-1.8	104	0.00
11 SURR	2-Methylnaphthalene-d10	0.642	0.638	0.6	100	0.00
12	2-Methylnaphthalene	0.768	0.771	-0.4	103	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
14 S	2,4,6-Tribromophenol	0.110	0.110	0.0	105	0.00
15 S	2-Fluorobiphenyl	1.410	1.442	-2.3	102	0.00
16	Acenaphthylene	2.112	2.101	0.5	101	0.00
17	Acenaphthene	1.356	1.351	0.4	99	0.00
18	Fluorene	1.738	1.736	0.1	101	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	96	0.00
20	4-Bromophenyl-phenylether	0.218	0.225	-3.2	101	0.00
21	Hexachlorobenzene	0.191	0.240	-25.7#	105	0.00
22	Pentachlorophenol	0.024	0.025	-4.2	97	0.00
23	Phenanthrene	1.141	1.221	-7.0	102	0.00
24	Anthracene	1.176	1.232	-4.8	99	0.00
25 SURR	Fluoranthene-d10	4.676	5.573	-19.2	100	0.00
26	Fluoranthene	1.420	1.729	-21.8	102	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	100	-0.01
28	Pvrene	1.399	1.419	-1.4	103	0.00
29 S	Terphenyl-d14	0.754	0.764	-1.3	103	0.00
30	Benzo(a)anthracene	1.302	1.350	-3.7	104	0.00
31	Chrysene	1.322	1.308	1.1	102	0.00
32	Bis(2-ethylhexyl)phthalate	3.193	3.089	3.3	106	-0.01
33	Indeno(1,2,3-cd)pyrene	1.222	1.248	-2.1	104	0.00
34 I	Perylene-d12	1.000	1.000	0.0	102	0.00
35	Benzo(b)fluoranthene	1.294	1.290	0.3	104	0.00
36	Benzo(k)fluoranthene	1.379	1.396	-1.2	104	0.00
37 C	Benzo(a)pyrene	1.255	1.244	0.9	101	0.00
38	Dibenzo(a,h)anthracene	1.127	1.132	-0.4	105	0.00
39	Benzo(a,h,i)perylene	1.055	1.076	-2.0	105	0.00

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