

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE042919\
 Data File : BE099414.D
 Acq On : 29 Apr 2019 18:38
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE042919

Quant Time: Apr 29 19:11:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE042919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 29 18:36:43 2019
 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	90	0.00
2	1,4-Dioxane	0.550	0.578	-5.1	110	0.00
3	n-Nitrosodimethylamine	0.332	0.347	-4.5	107	0.00
4 S	2-Fluorophenol	1.106	1.092	1.3	93	0.00
5 S	Phenol-d6	1.471	1.518	-3.2	95	0.00
6	bis(2-Chloroethyl)ether	1.269	1.252	1.3	91	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	90	0.00
8 S	Nitrobenzene-d5	0.375	0.383	-2.1	92	0.00
9	Naphthalene	4.359	4.481	-2.8	91	0.00
10	Hexachlorobutadiene	0.303	0.305	-0.7	89	0.00
11 SURR	2-Methylnaphthalene-d10	0.721	0.747	-3.6	94	0.00
12	2-Methylnaphthalene	0.758	0.755	0.4	90	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	92	0.00
14 S	2,4,6-Tribromophenol	0.252	0.244	3.2	92	0.00
15 S	2-Fluorobiphenyl	1.526	1.581	-3.6	93	0.00
16	Acenaphthylene	1.957	1.978	-1.1	92	0.00
17	Acenaphthene	1.118	1.110	0.7	91	0.00
18	Fluorene	1.649	1.686	-2.2	93	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
20	4-Bromophenyl-phenylether	0.233	0.227	2.6	92	0.00
21	Hexachlorobenzene	0.222	0.220	0.9	94	0.00
22	Pentachlorophenol	0.100	0.083	17.0	94	0.00
23	Phenanthrene	1.032	1.041	-0.9	95	0.00
24	Anthracene	0.984	0.980	0.4	95	0.00
25 SURR	Fluoranthene-d10	5.391	5.462	-1.3	96	0.00
26	Fluoranthene	1.546	1.579	-2.1	96	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	100	0.00
28	Pyrene	1.043	1.134	-8.7	96	0.00
29 S	Terphenyl-d14	0.733	0.799	-9.0	97	0.00
30	Benzo(a)anthracene	1.274	1.406	-10.4	98	0.00
31	Chrysene	1.247	1.418	-13.7	100	0.00
32	Bis(2-ethylhexyl)phthalate	2.216	2.149	3.0	89	0.00
33	Indeno(1,2,3-cd)pyrene	1.474	1.641	-11.3	98	0.00
34 I	Perylene-d12	1.000	1.000	0.0	100	0.00
35	Benzo(b)fluoranthene	1.153	1.209	-4.9	102	0.00
36	Benzo(k)fluoranthene	1.112	1.152	-3.6	98	0.00
37 C	Benzo(a)pyrene	1.090	1.124	-3.1	99	0.00
38	Dibenzo(a,h)anthracene	1.115	1.133	-1.6	98	0.01
39	Benzo(g,h,i)perylene	1.118	1.150	-2.9	99	0.00

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