

Data Path : Z:\HPCHEM1\BNA E\Data\BE041118\
 Data File : BE095974.D
 Acq On : 11 Apr 2018 17:09
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE041118

Quant Time: Apr 11 17:49:50 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE041118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 11 17:26:43 2018
 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	108	0.00
2	1,4-Dioxane	0.618	0.575	7.0	108	0.00
3	n-Nitrosodimethylamine	0.757	0.714	5.7	111	0.00
4 S	2-Fluorophenol	1.237	1.154	6.7	106	0.00
5 S	Phenol-d6	1.707	1.589	6.9	106	0.00
6	bis(2-Chloroethyl)ether	1.502	1.382	8.0	104	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
8 S	Nitrobenzene-d5	0.469	0.440	6.2	109	0.00
9	Naphthalene	4.217	3.939	6.6	106	0.00
10	Hexachlorobutadiene	0.187	0.220	-17.6	139	0.00
11 SURR	2-Methylnaphthalene-d10	0.634	0.597	5.8	108	0.00
12	2-Methylnaphthalene	0.698	0.651	6.7	108	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
14 S	2,4,6-Tribromophenol	0.193	0.176	8.8	109	0.00
15 S	2-Fluorobiphenyl	1.711	1.619	5.4	108	0.00
16	Acenaphthylene	2.136	1.956	8.4	108	0.01
17	Acenaphthene	1.284	1.206	6.1	111	0.00
18	Fluorene	1.588	1.498	5.7	110	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	109	0.01
20	4-Bromophenyl-phenylether	0.238	0.221	7.1	108	0.00
21	Hexachlorobenzene	0.237	0.225	5.1	107	0.00
22	Pentachlorophenol	0.078	0.065	16.7	112	0.00
23	Phenanthrene	1.124	1.064	5.3	109	0.00
24	Anthracene	1.074	1.006	6.3	111	0.00
25 SURR	Fluoranthene-d10	5.430	5.057	6.9	110	0.00
26	Fluoranthene	1.477	1.378	6.7	110	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	107	0.00
28	Pvrene	0.788	0.778	1.3	118	0.00
29 S	Terphenyl-d14	0.884	0.840	5.0	111	0.00
30	Benzo(a)anthracene	1.283	1.174	8.5	107	0.00
31	Chrysene	1.239	1.168	5.7	107	0.00
32	Bis(2-ethylhexyl)phthalate	3.311	2.983	9.9	104	0.00
33	Indeno(1,2,3-cd)pyrene	1.201	1.102	8.2	105	0.00
34 I	Perylene-d12	1.000	1.000	0.0	104	0.00
35	Benzo(b)fluoranthene	1.200	1.132	5.7	103	0.00
36	Benzo(k)fluoranthene	1.234	1.115	9.6	101	0.00
37 C	Benzo(a)pyrene	1.153	1.073	6.9	103	0.00
38	Dibenzo(a,h)anthracene	1.017	0.929	8.7	103	0.00
39	Benzo(a,h,i)perylene	1.076	1.012	5.9	104	0.00

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