

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110518\
 Data File : BE098153.D
 Acq On : 5 Nov 2018 9:41
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 06 10:21:34 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 13:00:52 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	81	-0.01
2	1,4-Dioxane	0.596	0.575	3.5	78	0.00
3	n-Nitrosodimethylamine	0.797	0.784	1.6	85	0.00
4 S	2-Fluorophenol	1.187	1.134	4.5	76	0.00
5 S	Phenol-d6	1.955	1.719	12.1	71	0.00
6	bis(2-Chloroethyl)ether	1.431	1.406	1.7	82	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	78	0.00
8 S	Nitrobenzene-d5	0.429	0.380	11.4	71	0.00
9	Naphthalene	3.998	4.208	-5.3	84	0.00
10	Hexachlorobutadiene	0.262	0.274	-4.6	84	0.00
11 SURR	2-Methylnaphthalene-d10	0.705	0.704	0.1	78	0.00
12	2-Methylnaphthalene	0.714	0.742	-3.9	83	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	72	-0.01
14 S	2,4,6-Tribromophenol	0.189	0.166	12.2	75	0.00
15 S	2-Fluorobiphenyl	1.645	1.696	-3.1	82	0.00
16	Acenaphthylene	1.825	1.892	-3.7	80	0.00
17	Acenaphthene	1.096	1.153	-5.2	79	0.00
18	Fluorene	1.454	1.495	-2.8	80	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	76	0.00
20	4-Bromophenyl-phenylether	0.244	0.255	-4.5	82	0.00
21	Hexachlorobenzene	0.255	0.271	-6.3	85	0.00
22	Pentachlorophenol	0.057	0.041	28.1#	62	0.00
23	Phenanthrene	0.991	1.034	-4.3	83	0.00
24	Anthracene	0.942	0.959	-1.8	83	-0.01
25 SURR	Fluoranthene-d10	5.159	5.388	-4.4	85	0.00
26	Fluoranthene	1.241	1.316	-6.0	86	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	78	0.00
28	Pyrene	1.092	1.122	-2.7	87	0.00
29 S	Terphenyl-d14	0.930	0.959	-3.1	88	0.00
30	Benzo(a)anthracene	1.169	1.189	-1.7	86	0.00
31	Chrysene	1.156	1.205	-4.2	88	0.00
32	Bis(2-ethylhexyl)phthalate	4.140	1.834	55.7#	43#	0.00
33	Indeno(1,2,3-cd)pyrene	1.203	1.198	0.4	84	0.00
34 I	Perylene-d12	1.000	1.000	0.0	80	0.00
35	Benzo(b)fluoranthene	1.127	1.144	-1.5	86	0.00
36	Benzo(k)fluoranthene	1.166	1.158	0.7	85	0.00
37 C	Benzo(a)pyrene	1.101	1.084	1.5	83	0.00
38	Dibenzo(a,h)anthracene	1.027	1.022	0.5	84	0.00
39	Benzo(g,h,i)perylene	1.022	1.013	0.9	84	-0.01

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