

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011719\
 Data File : BE098646.D
 Acq On : 17 Jan 2019 9:18
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 17 12:40:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE011519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 15 14:30:55 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	84	0.00
2	1,4-Dioxane	0.789	0.724	8.2	89	0.00
3	n-Nitrosodimethylamine	0.579	0.559	3.5	85	0.01
4 S	2-Fluorophenol	1.067	1.045	2.1	84	0.00
5 S	Phenol-d6	1.519	1.504	1.0	83	0.00
6	bis(2-Chloroethyl)ether	1.149	0.992	13.7	78	-0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	80	0.00
8 S	Nitrobenzene-d5	0.317	0.337	-6.3	87	0.00
9	Naphthalene	3.984	3.822	4.1	81	0.00
10	Hexachlorobutadiene	0.285	0.294	-3.2	86	0.00
11 SURR	2-Methylnaphthalene-d10	0.658	0.664	-0.9	87	0.00
12	2-Methylnaphthalene	0.660	0.649	1.7	85	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	86	0.00
14 S	2,4,6-Tribromophenol	0.174	0.167	4.0	84	0.00
15 S	2-Fluorobiphenyl	1.638	1.458	11.0	84	0.00
16	Acenaphthylene	1.732	1.564	9.7	83	0.00
17	Acenaphthene	1.087	1.003	7.7	83	0.00
18	Fluorene	1.414	1.300	8.1	82	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	84	0.00
20	4-Bromophenyl-phenylether	0.212	0.202	4.7	86	0.00
21	Hexachlorobenzene	0.256	0.244	4.7	84	0.00
22	Pentachlorophenol	0.075	0.078	-4.0	111	0.00
23	Phenanthrene	0.948	0.864	8.9	82	0.00
24	Anthracene	0.812	0.728	10.3	83	0.00
25 SURR	Fluoranthene-d10	4.955	4.685	5.4	86	0.00
26	Fluoranthene	1.220	1.136	6.9	85	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	85	0.00
28	Pvrene	1.101	1.034	6.1	84	0.00
29 S	Terphenyl-d14	0.885	0.822	7.1	86	0.00
30	Benzo(a)anthracene	1.075	1.042	3.1	88	0.00
31	Chrysene	1.115	1.040	6.7	84	0.00
32	Bis(2-ethylhexyl)phthalate	2.281	2.049	10.2	86	0.00
33	Indeno(1,2,3-cd)pyrene	1.150	1.098	4.5	85	0.00
34 I	Perylene-d12	1.000	1.000	0.0	87	0.00
35	Benzo(b)fluoranthene	1.091	1.019	6.6	87	0.00
36	Benzo(k)fluoranthene	1.193	1.121	6.0	84	0.00
37 C	Benzo(a)pyrene	1.072	1.003	6.4	86	0.00
38	Dibenzo(a,h)anthracene	0.949	0.907	4.4	88	-0.01
39	Benzo(a,h,i)perylene	1.040	0.980	5.8	85	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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