

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE113018\
 Data File : BE098283.D
 Acq On : 30 Nov 2018 9:36
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 30 11:04:44 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE112918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 29 13:50:43 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	86	0.00
2	1,4-Dioxane	0.742	0.722	2.7	93	0.00
3	n-Nitrosodimethylamine	0.768	0.665	13.4	76	0.00
4 S	2-Fluorophenol	1.068	1.087	-1.8	89	0.01
5 S	Phenol-d6	1.505	1.415	6.0	84	0.00
6	bis(2-Chloroethyl)ether	1.385	1.328	4.1	84	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
8 S	Nitrobenzene-d5	0.405	0.363	10.4	81	0.00
9	Naphthalene	4.019	3.989	0.7	88	-0.01
10	Hexachlorobutadiene	0.258	0.258	0.0	89	0.00
11 SURR	2-Methylnaphthalene-d10	0.650	0.657	-1.1	87	0.00
12	2-Methylnaphthalene	0.667	0.645	3.3	86	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	86	0.00
14 S	2,4,6-Tribromophenol	0.190	0.169	11.1	87	0.00
15 S	2-Fluorobiphenyl	1.657	1.574	5.0	87	0.00
16	Acenaphthylene	1.867	1.848	1.0	88	0.00
17	Acenaphthene	1.113	1.088	2.2	86	0.00
18	Fluorene	1.475	1.486	-0.7	87	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
20	4-Bromophenyl-phenylether	0.226	0.207	8.4	83	0.00
21	Hexachlorobenzene	0.234	0.225	3.8	86	0.00
22	Pentachlorophenol	0.076	0.066	13.2	98	0.00
23	Phenanthrene	0.993	0.959	3.4	87	0.00
24	Anthracene	0.925	0.866	6.4	86	0.00
25 SURR	Fluoranthene-d10	5.397	5.458	-1.1	93	0.00
26	Fluoranthene	1.356	1.383	-2.0	95	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	113	0.00
28	Pvrene	1.108	1.027	7.3	98	0.00
29 S	Terphenyl-d14	0.900	0.799	11.2	97	0.00
30	Benzo(a)anthracene	1.178	1.117	5.2	111	0.00
31	Chrysene	1.196	1.183	1.1	114	-0.01
32	Bis(2-ethylhexyl)phthalate	2.821	2.433	13.8	100	0.00
33	Indeno(1,2,3-cd)pyrene	1.210	1.220	-0.8	122	0.00
34 I	Perylene-d12	1.000	1.000	0.0	121	0.00
35	Benzo(b)fluoranthene	1.173	1.083	7.7	114	0.00
36	Benzo(k)fluoranthene	1.231	1.224	0.6	125	0.00
37 C	Benzo(a)pyrene	1.130	1.102	2.5	121	0.00
38	Dibenzo(a,h)anthracene	1.071	1.046	2.3	121	0.00
39	Benzo(a,h,i)perylene	1.073	1.033	3.7	121	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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