

Data Path : U:\HPCHEM1\BNA E\Data\BE082217\
 Data File : BE093871.D
 Acq On : 22 Aug 2017 12:18
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 23 01:20:14 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 07 14:30:21 2017
 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	78	0.00
2	1,4-Dioxane	0.603	0.463	23.2	59	0.00
3	n-Nitrosodimethylamine	0.625	0.541	13.4	66	0.01
4 S	2-Fluorophenol	1.294	1.112	14.1	68	0.00
5 S	Phenol-d6	1.976	1.672	15.4	68	0.01
6	bis(2-Chloroethyl)ether	1.530	1.310	14.4	69	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	79	0.00
8 S	Nitrobenzene-d5	0.342	0.302	11.7	73	0.00
9	Naphthalene	4.587	4.036	12.0	70	0.00
10	Hexachlorobutadiene	0.170	0.141	17.1	66	0.00
11 SURR	2-Methylnaphthalene-d10	0.643	0.564	12.3	69	0.00
12	2-Methylnaphthalene	0.774	0.669	13.6	69	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	77	0.00
14 S	2,4,6-Tribromophenol	0.110	0.088	20.0	68	0.03
15 S	2-Fluorobiphenyl	1.577	1.379	12.6	67	0.00
16	Acenaphthylene	2.011	1.750	13.0	69	0.00
17	Acenaphthene	1.355	1.152	15.0	66	0.00
18	Fluorene	1.752	1.507	14.0	67	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	71	0.00
20	4-Bromophenyl-phenylether	0.113	0.130	-15.0	83	0.00
21	Hexachlorobenzene	0.121	0.124	-2.5	70	0.00
22	Pentachlorophenol	0.084	0.055	34.5#	49#	0.00
23	Phenanthrene	0.852	0.808	5.2	68	0.00
24	Anthracene	1.294	1.114	13.9	62	0.00
25 SURR	Fluoranthene-d10	4.564	4.262	6.6	62	0.00
26	Fluoranthene	1.397	1.269	9.2	61	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	68	0.01
28	Pvrene	1.310	1.144	12.7	61	0.00
29 S	Terphenyl-d14	0.713	0.620	13.0	61	0.00
30	Benzo(a)anthracene	1.269	1.052	17.1	58	0.00
31	Chrysene	1.295	1.156	10.7	62	0.00
32	Bis(2-ethylhexyl)phthalate	2.721	2.432	10.6	60	0.00
33	Indeno(1,2,3-cd)pyrene	1.248	0.893	28.4#	49#	0.04
34 I	Perylene-d12	1.000	1.000	0.0	67	0.01
35	Benzo(b)fluoranthene	1.387	1.074	22.6	51	0.01
36	Benzo(k)fluoranthene	1.386	1.240	10.5	62	0.01
37 C	Benzo(a)pyrene	1.309	1.097	16.2	56	0.02
38	Dibenzo(a,h)anthracene	1.225	0.896	26.9#	50#	0.04
39	Benzo(a,h,i)perylene	1.229	0.914	25.6#	51	0.04

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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