

Data Path : Z:\HPCHEM1\BNA E\Data\BE032318\
 Data File : BE095830.D
 Acq On : 23 Mar 2018 19:25
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 23 23:23:58 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE032318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 23 12:41:11 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	107	0.00
2	1,4-Dioxane	0.655	0.600	8.4	103	0.00
3	n-Nitrosodimethylamine	0.961	0.812	15.5	87	0.00
4 S	2-Fluorophenol	1.319	1.167	11.5	91	0.00
5 S	Phenol-d6	1.813	1.666	8.1	94	0.00
6	bis(2-Chloroethyl)ether	1.564	1.449	7.4	95	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
8 S	Nitrobenzene-d5	0.473	0.430	9.1	96	0.00
9	Naphthalene	4.569	4.151	9.1	93	0.00
10	Hexachlorobutadiene	0.314	0.260	17.2	84	0.00
11 SURR	2-Methylnaphthalene-d10	0.668	0.612	8.4	94	0.00
12	2-Methylnaphthalene	0.780	0.728	6.7	96	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
14 S	2,4,6-Tribromophenol	0.231	0.200	13.4	98	0.00
15 S	2-Fluorobiphenyl	1.686	1.499	11.1	95	-0.01
16	Acenaphthylene	2.288	1.998	12.7	95	0.00
17	Acenaphthene	1.395	1.222	12.4	96	0.00
18	Fluorene	1.835	1.607	12.4	95	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
20	4-Bromophenyl-phenylether	0.240	0.215	10.4	98	0.00
21	Hexachlorobenzene	0.240	0.213	11.2	97	0.00
22	Pentachlorophenol	0.069	0.043	37.7#	82	0.00
23	Phenanthrene	1.197	1.091	8.9	98	0.00
24	Anthracene	1.197	1.085	9.4	100	0.00
25 SURR	Fluoranthene-d10	5.940	5.442	8.4	100	0.00
26	Fluoranthene	1.663	1.531	7.9	101	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	111	-0.01
28	Pvrene	1.629	1.548	5.0	96	-0.01
29 S	Terphenyl-d14	0.850	0.764	10.1	99	0.00
30	Benzo(a)anthracene	1.387	1.252	9.7	98	0.00
31	Chrysene	1.266	1.184	6.5	103	0.00
32	Bis(2-ethylhexyl)phthalate	4.242	3.425	19.3	96	0.00
33	Indeno(1,2,3-cd)pyrene	1.364	1.252	8.2	101	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	113	0.00
35	Benzo(b)fluoranthene	1.323	1.193	9.8	101	0.00
36	Benzo(k)fluoranthene	1.324	1.206	8.9	102	0.00
37 C	Benzo(a)pyrene	1.240	1.132	8.7	101	0.00
38	Dibenzo(a,h)anthracene	1.178	1.060	10.0	99	-0.01
39	Benzo(a,h,i)perylene	1.184	1.059	10.6	99	-0.01

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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