

Data Path : Z:\HPCHEM1\BNA E\DATA\BE051917\
 Data File : BE093063.D
 Acq On : 19 May 2017 16:29
 Operator : SJ/MA
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: May 19 18:06:36 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE051517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 15 13:35:57 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	80	0.00
2	1,4-Dioxane	0.810	0.729	10.0	85	0.00
3	n-Nitrosodimethylamine	0.724	0.648	10.5	82	0.00
4 S	2-Fluorophenol	1.226	1.093	10.8	80	0.00
5 S	Phenol-d6	1.656	1.462	11.7	81	0.00
6	bis(2-Chloroethyl)ether	1.349	1.190	11.8	81	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	82	0.00
8 S	Nitrobenzene-d5	0.350	0.325	7.1	88	0.00
9	Naphthalene	1.054	0.928	12.0	81	-0.02
10	Hexachlorobutadiene	0.138	0.128	7.2	86	0.00
11 SURR	2-Methylnaphthalene-d10	0.586	0.515	12.1	81	0.00
12	2-Methylnaphthalene	0.653	0.572	12.4	81	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	82	0.00
14 S	2,4,6-Tribromophenol	0.118	0.113	4.2	91	-0.01
15 S	2-Fluorobiphenyl	1.645	1.447	12.0	80	0.00
16	Acenaphthylene	7.015	6.070	13.5	84	0.00
17	Acenaphthene	1.263	1.111	12.0	82	0.00
18	Fluorene	1.545	1.366	11.6	82	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
20	4-Bromophenyl-phenylether	0.124	0.111	10.5	97	0.00
21	Hexachlorobenzene	0.163	0.110	32.5#	73	0.00
22	Pentachlorophenol	0.050	0.037	26.0#	97	0.00
23	Phenanthrene	0.867	0.670	22.7	86	0.00
24	Anthracene	0.816	0.569	30.3#	83	0.00
25 SURR	Fluoranthene-d10	1.051	0.834	20.6	85	0.00
26	Fluoranthene	4.922	3.986	19.0	89	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	88	-0.02
28	Pvrene	4.819	4.211	12.6	84	-0.02
29 S	Terphenyl-d14	0.747	0.592	20.7	76	0.00
30	Benzo(a)anthracene	1.088	1.016	6.6	92	0.00
31	Chrysene	1.163	0.975	16.2	83	0.00
32	Bis(2-ethylhexyl)phthalate	0.627	0.463	26.2#	93	0.00
33	Indeno(1,2,3-cd)pyrene	4.256	3.881	8.8	91	0.00
34 I	Perylene-d12	1.000	1.000	0.0	87	0.00
35	Benzo(b)fluoranthene	1.264	1.194	5.5	96	0.00
36	Benzo(k)fluoranthene	1.268	1.014	20.0	79	0.00
37 C	Benzo(a)pyrene	1.157	1.000	13.6	88	0.00
38	Dibenzo(a,h)anthracene	1.076	0.990	8.0	93	0.00
39	Benzo(a,h,i)perylene	4.686	4.226	9.8	90	-0.02

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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