

Data Path : Z:\HPCHEM1\BNA_E\Data\BE070117\
 Data File : BE093386.D
 Acq On : 1 Jul 2017 11:08
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 03 05:21:10 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 01 15:43:24 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	93	-0.01
2	1,4-Dioxane	0.826	0.800	3.1	93	0.00
3	n-Nitrosodimethylamine	0.497	0.407	18.1	91	-0.01
4 S	2-Fluorophenol	1.247	1.203	3.5	93	0.00
5 S	Phenol-d6	1.861	1.702	8.5	91	0.00
6	bis(2-Chloroethyl)ether	1.312	1.299	1.0	100	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
8 S	Nitrobenzene-d5	0.238	0.229	3.8	102	-0.02
9	Naphthalene	4.155	4.081	1.8	94	-0.02
10	Hexachlorobutadiene	0.167	0.171	-2.4	95	0.00
11 SURR	2-Methylnaphthalene-d10	0.639	0.611	4.4	92	0.00
12	2-Methylnaphthalene	0.708	0.676	4.5	91	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	88	-0.02
14 S	2,4,6-Tribromophenol	0.172	0.150	12.8	84	0.00
15 S	2-Fluorobiphenyl	1.566	1.623	-3.6	92	0.00
16	Acenaphthylene	1.739	1.616	7.1	86	-0.02
17	Acenaphthene	1.224	1.202	1.8	89	0.00
18	Fluorene	1.683	1.702	-1.1	90	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	85	0.00
20	4-Bromophenyl-phenylether	0.091	0.126	-38.5#	116	0.00
21	Hexachlorobenzene	0.153	0.183	-19.6	101	0.00
22	Pentachlorophenol	0.018	0.016	11.1	92	-0.03
23	Phenanthrene	0.968	1.100	-13.6	93	0.00
24	Anthracene	0.974	0.826	15.2	78	0.00
25 SURR	Fluoranthene-d10	3.688	3.658	0.8	89	0.00
26	Fluoranthene	1.049	1.039	1.0	89	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	94	-0.01
28	Pyrene	1.056	0.988	6.4	90	0.00
29 S	Terphenyl-d14	0.742	0.720	3.0	93	0.00
30	Benzo(a)anthracene	1.088	1.028	5.5	93	0.00
31	Chrysene	1.137	1.122	1.3	94	0.00
32	Bis(2-ethylhexyl)phthalate	1.612	1.318	18.2	92	-0.01
33	Indeno(1,2,3-cd)pyrene	1.112	1.120	-0.7	98	-0.02
34 I	Perylene-d12	1.000	1.000	0.0	97	0.00
35	Benzo(b)fluoranthene	1.256	1.214	3.3	98	0.00
36	Benzo(k)fluoranthene	1.260	1.217	3.4	98	-0.01
37 C	Benzo(a)pyrene	1.116	1.059	5.1	97	0.00
38	Dibenzo(a,h)anthracene	1.137	1.097	3.5	98	0.00
39	Benzo(g,h,i)perylene	1.159	1.131	2.4	98	0.00

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