

Data Path : Z:\HPCHEM1\BNA E\Data\BE032318\
 Data File : BE095821.D
 Acq On : 23 Mar 2018 12:21
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE032318

Quant Time: Mar 23 12:57:01 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	101	-0.01
2	1,4-Dioxane	0.655	0.661	-0.9	108	0.00
3	n-Nitrosodimethylamine	0.961	0.879	8.5	89	0.00
4 S	2-Fluorophenol	1.319	1.275	3.3	94	-0.01
5 S	Phenol-d6	1.813	1.725	4.9	92	0.00
6	bis(2-Chloroethyl)ether	1.564	1.458	6.8	91	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	96	0.00
8 S	Nitrobenzene-d5	0.473	0.472	0.2	95	0.00
9	Naphthalene	4.569	4.504	1.4	92	0.00
10	Hexachlorobutadiene	0.314	0.313	0.3	92	0.00
11 SURR	2-Methylnaphthalene-d10	0.668	0.679	-1.6	94	0.00
12	2-Methylnaphthalene	0.780	0.794	-1.8	95	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
14 S	2,4,6-Tribromophenol	0.231	0.216	6.5	95	0.00
15 S	2-Fluorobiphenyl	1.686	1.648	2.3	95	0.00
16	Acenaphthylene	2.288	2.248	1.7	97	0.00
17	Acenaphthene	1.395	1.368	1.9	97	0.00
18	Fluorene	1.835	1.800	1.9	96	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
20	4-Bromophenyl-phenylether	0.240	0.237	1.3	97	0.00
21	Hexachlorobenzene	0.240	0.235	2.1	96	0.00
22	Pentachlorophenol	0.069	0.051	26.1#	88	0.00
23	Phenanthrene	1.197	1.197	0.0	97	0.00
24	Anthracene	1.197	1.179	1.5	97	0.00
25 SURR	Fluoranthene-d10	5.940	5.957	-0.3	98	0.00
26	Fluoranthene	1.663	1.667	-0.2	99	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	100	0.00
28	Pvrene	1.629	1.799	-10.4	100	0.00
29 S	Terphenyl-d14	0.850	0.847	0.4	98	0.00
30	Benzo(a)anthracene	1.387	1.405	-1.3	99	0.00
31	Chrysene	1.266	1.278	-0.9	99	0.00
32	Bis(2-ethylhexyl)phthalate	4.242	3.803	10.3	95	0.00
33	Indeno(1,2,3-cd)pyrene	1.364	1.388	-1.8	100	0.00
34 I	Perylene-d12	1.000	1.000	0.0	101	0.00
35	Benzo(b)fluoranthene	1.323	1.338	-1.1	101	0.00
36	Benzo(k)fluoranthene	1.324	1.350	-2.0	102	0.00
37 C	Benzo(a)pyrene	1.240	1.247	-0.6	100	0.00
38	Dibenzo(a,h)anthracene	1.178	1.198	-1.7	100	0.00
39	Benzo(a,h,i)perylene	1.184	1.162	1.9	98	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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