

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE050719\
 Data File : BE099565.D
 Acq On : 7 May 2019 14:30
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICSVBE050719

Quant Time: May 07 15:21:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE050719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 07 14:10:11 2019
 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	93	0.02
2	1,4-Dioxane	0.565	0.567	-0.4	113	0.00
3	n-Nitrosodimethylamine	0.374	0.349	6.7	96	0.00
4 S	2-Fluorophenol	1.109	1.125	-1.4	105	0.00
5 S	Phenol-d6	1.459	1.428	2.1	99	0.02
6	bis(2-Chloroethyl)ether	1.216	1.209	0.6	103	0.02
7 I	Naphthalene-d8	1.000	1.000	0.0	91	0.01
8 S	Nitrobenzene-d5	0.385	0.381	1.0	100	0.03
9	Naphthalene	4.237	4.377	-3.3	104	0.01
10	Hexachlorobutadiene	0.308	0.324	-5.2	106	0.01
11 SURR	2-Methylnaphthalene-d10	0.697	0.729	-4.6	106	0.03
12	2-Methylnaphthalene	0.718	0.731	-1.8	101	0.01
13 I	Acenaphthene-d10	1.000	1.000	0.0	89	0.01
14 S	2,4,6-Tribromophenol	0.266	0.250	6.0	98	0.01
15 S	2-Fluorobiphenyl	1.516	1.523	-0.5	103	0.01
16	Acenaphthylene	1.930	1.889	2.1	99	0.01
17	Acenaphthene	1.078	1.049	2.7	97	0.01
18	Fluorene	1.540	1.534	0.4	99	0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	91	0.01
20	4-Bromophenyl-phenylether	0.227	0.217	4.4	99	0.01
21	Hexachlorobenzene	0.226	0.226	0.0	101	0.00
22	Pentachlorophenol	0.093	0.081	12.9	99	0.01
23	Phenanthrene	0.986	1.004	-1.8	103	0.01
24	Anthracene	0.948	0.947	0.1	104	0.01
25 SURR	Fluoranthene-d10	5.386	5.358	0.5	102	0.00
26	Fluoranthene	1.518	1.512	0.4	101	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	101	0.00
28	Pvrene	1.072	1.069	0.3	102	0.00
29 S	Terphenyl-d14	0.741	0.741	0.0	103	0.01
30	Benzo(a)anthracene	1.239	1.192	3.8	99	0.00
31	Chrysene	1.223	1.234	-0.9	102	0.01
32	Bis(2-ethylhexyl)phthalate	1.964	1.943	1.1	103	0.00
33	Indeno(1,2,3-cd)pyrene	1.452	1.415	2.5	100	0.01
34 I	Perylene-d12	1.000	1.000	0.0	91	0.01
35	Benzo(b)fluoranthene	1.103	1.122	-1.7	101	0.00
36	Benzo(k)fluoranthene	1.086	1.125	-3.6	104	0.00
37 C	Benzo(a)pyrene	1.064	1.076	-1.1	101	0.00
38	Dibenzo(a,h)anthracene	1.100	1.107	-0.6	101	0.02
39	Benzo(a,h,i)perylene	1.102	1.114	-1.1	100	0.01

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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