

Data Path : Z:\HPCHEM1\BNA E\DATA\BE051117\
 Data File : BE092995.D
 Acq On : 11 May 2017 14:27
 Operator : SJ/MA
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE051117

Quant Time: May 11 16:40:41 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE051117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 11 14:48:08 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	88	-0.02
2	1,4-Dioxane	0.802	0.781	2.6	101	0.00
3	n-Nitrosodimethylamine	0.688	0.676	1.7	100	0.00
4 S	2-Fluorophenol	1.097	1.061	3.3	99	0.00
5 S	Phenol-d6	1.376	1.359	1.2	101	-0.02
6	bis(2-Chloroethyl)ether	1.347	1.353	-0.4	102	-0.02
7 I	Naphthalene-d8	1.000	1.000	0.0	89	0.00
8 S	Nitrobenzene-d5	0.313	0.305	2.6	100	-0.02
9	Naphthalene	1.065	1.078	-1.2	100	0.00
10	Hexachlorobutadiene	0.146	0.149	-2.1	102	-0.02
11 SURR	2-Methylnaphthalene-d10	0.567	0.571	-0.7	102	-0.02
12	2-Methylnaphthalene	0.624	0.617	1.1	100	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	90	0.00
14 S	2,4,6-Tribromophenol	0.109	0.102	6.4	104	0.00
15 S	2-Fluorobiphenyl	1.585	1.567	1.1	101	0.00
16	Acenaphthylene	6.788	6.382	6.0	102	0.00
17	Acenaphthene	1.236	1.218	1.5	101	0.00
18	Fluorene	1.540	1.505	2.3	100	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	93	0.00
20	4-Bromophenyl-phenylether	0.128	0.123	3.9	101	0.00
21	Hexachlorobenzene	0.191	0.200	-4.7	105	0.00
22	Pentachlorophenol	0.048	0.045	6.3	101	0.00
23	Phenanthrene	0.959	0.917	4.4	104	0.01
24	Anthracene	0.828	0.877	-5.9	107	0.00
25 SURR	Fluoranthene-d10	1.059	1.040	1.8	102	-0.02
26	Fluoranthene	5.138	5.156	-0.4	104	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	94	0.00
28	Pvrene	5.018	5.163	-2.9	107	0.00
29 S	Terphenyl-d14	0.710	0.680	4.2	108	0.00
30	Benzo(a)anthracene	0.983	0.912	7.2	108	0.00
31	Chrysene	1.228	1.276	-3.9	109	0.00
32	Bis(2-ethylhexyl)phthalate	0.434	0.353	18.7	107	0.00
33	Indeno(1,2,3-cd)pyrene	3.923	4.196	-7.0	114	0.00
34 I	Perylene-d12	1.000	1.000	0.0	95	0.00
35	Benzo(b)fluoranthene	1.290	1.202	6.8	96	0.00
36	Benzo(k)fluoranthene	1.289	1.358	-5.4	125	0.00
37 C	Benzo(a)pyrene	1.122	1.060	5.5	106	0.00
38	Dibenzo(a,h)anthracene	1.016	1.083	-6.6	119	0.02
39	Benzo(a,h,i)perylene	4.629	4.818	-4.1	109	0.00

Data Path : Z:\HPCHEM1\BNA E\DATA\BE051117\
Data File : BE092995.D
Acq On : 11 May 2017 14:27
Operator : SJ/MA
Sample : SSTDICV0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
ICVBE051117

Quant Time: May 11 16:40:41 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE051117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 11 14:48:08 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)

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