

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE092718\
 Data File : BE097696.D
 Acq On : 27 Sep 2018 16:17
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE092718

Quant Time: Sep 27 17:01:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE092718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 27 15:45:00 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	94	0.00
2	1,4-Dioxane	0.834	0.779	6.6	95	0.00
3	n-Nitrosodimethylamine	0.501	0.438	12.6	87	0.00
4 S	2-Fluorophenol	0.989	0.852	13.9	87	0.00
5 S	Phenol-d6	1.363	1.190	12.7	90	0.01
6	bis(2-Chloroethyl)ether	1.335	1.274	4.6	95	0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
8 S	Nitrobenzene-d5	0.284	0.256	9.9	90	0.00
9	Naphthalene	3.951	3.806	3.7	97	0.00
10	Hexachlorobutadiene	0.185	0.180	2.7	100	0.00
11 SURR	2-Methylnaphthalene-d10	0.622	0.588	5.5	97	0.00
12	2-Methylnaphthalene	0.604	0.567	6.1	96	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
14 S	2,4,6-Tribromophenol	0.156	0.126	19.2	91	0.00
15 S	2-Fluorobiphenyl	1.680	1.626	3.2	98	0.00
16	Acenaphthylene	1.707	1.601	6.2	97	0.00
17	Acenaphthene	1.149	1.140	0.8	98	0.00
18	Fluorene	1.468	1.408	4.1	95	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
20	4-Bromophenyl-phenylether	0.204	0.192	5.9	96	0.00
21	Hexachlorobenzene	0.229	0.222	3.1	100	0.00
22	Pentachlorophenol	0.053	0.044	17.0	85	0.01
23	Phenanthrene	1.034	0.981	5.1	97	0.00
24	Anthracene	0.851	0.753	11.5	94	0.00
25 SURR	Fluoranthene-d10	4.850	4.594	5.3	98	0.00
26	Fluoranthene	1.284	1.218	5.1	98	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	97	0.00
28	Pvrene	1.041	1.014	2.6	97	0.00
29 S	Terphenyl-d14	0.864	0.837	3.1	98	0.00
30	Benzo(a)anthracene	0.989	0.889	10.1	93	0.00
31	Chrysene	1.238	1.262	-1.9	102	0.00
32	Bis(2-ethylhexyl)phthalate	1.220	0.986	19.2	89	0.00
33	Indeno(1,2,3-cd)pyrene	1.307	1.230	5.9	96	0.00
34 I	Perylene-d12	1.000	1.000	0.0	96	0.00
35	Benzo(b)fluoranthene	1.156	1.117	3.4	99	0.00
36	Benzo(k)fluoranthene	1.334	1.294	3.0	96	0.00
37 C	Benzo(a)pyrene	1.109	1.023	7.8	95	0.00
38	Dibenzo(a,h)anthracene	1.203	1.118	7.1	96	0.00
39	Benzo(a,h,i)perylene	1.220	1.154	5.4	96	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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