

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE050919\
 Data File : BE099589.D
 Acq On : 9 May 2019 13:41
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE050919

Quant Time: May 09 16:06:11 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE050919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 09 13:32:58 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	80	0.01
2	1,4-Dioxane	0.564	0.614	-8.9	103	0.00
3	n-Nitrosodimethylamine	0.349	0.337	3.4	93	0.00
4 S	2-Fluorophenol	1.094	1.107	-1.2	94	0.00
5 S	Phenol-d6	1.437	1.375	4.3	88	0.01
6	bis(2-Chloroethyl)ether	1.195	1.247	-4.4	96	0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	80	0.01
8 S	Nitrobenzene-d5	0.375	0.362	3.5	87	0.01
9	Naphthalene	4.262	4.319	-1.3	91	0.00
10	Hexachlorobutadiene	0.308	0.306	0.6	88	0.01
11 SURR	2-Methylnaphthalene-d10	0.670	0.671	-0.1	88	0.01
12	2-Methylnaphthalene	0.696	0.706	-1.4	90	0.01
13 I	Acenaphthene-d10	1.000	1.000	0.0	78	0.00
14 S	2,4,6-Tribromophenol	0.237	0.243	-2.5	92	0.01
15 S	2-Fluorobiphenyl	1.551	1.589	-2.5	89	0.00
16	Acenaphthylene	1.936	1.923	0.7	87	0.01
17	Acenaphthene	1.081	1.087	-0.6	87	0.00
18	Fluorene	1.543	1.556	-0.8	89	0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	78	0.00
20	4-Bromophenyl-phenylether	0.228	0.220	3.5	87	0.01
21	Hexachlorobenzene	0.229	0.232	-1.3	88	0.01
22	Pentachlorophenol	0.092	0.070	23.9	81	0.01
23	Phenanthrene	0.993	1.006	-1.3	89	0.00
24	Anthracene	0.932	0.914	1.9	89	0.00
25 SURR	Fluoranthene-d10	5.261	5.311	-1.0	89	0.00
26	Fluoranthene	1.494	1.525	-2.1	89	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	85	0.00
28	Pvrene	1.054	1.075	-2.0	87	0.00
29 S	Terphenyl-d14	0.711	0.720	-1.3	87	0.00
30	Benzo(a)anthracene	1.221	1.226	-0.4	86	0.01
31	Chrysene	1.221	1.285	-5.2	90	0.00
32	Bis(2-ethylhexyl)phthalate	1.755	1.735	1.1	87	0.01
33	Indeno(1,2,3-cd)pyrene	1.524	1.600	-5.0	92	0.01
34 I	Perylene-d12	1.000	1.000	0.0	81	0.00
35	Benzo(b)fluoranthene	1.105	1.131	-2.4	91	0.00
36	Benzo(k)fluoranthene	1.098	1.082	1.5	87	0.00
37 C	Benzo(a)pyrene	1.071	1.071	0.0	90	0.00
38	Dibenzo(a,h)anthracene	1.119	1.125	-0.5	91	0.01
39	Benzo(a,h,i)perylene	1.133	1.157	-2.1	91	0.01

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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