

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE100319\
 Data File : BE100601.D
 Acq On : 3 Oct 2019 21:04
 Operator : JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE100319

Quant Time: Oct 04 13:08:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE100319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 04 13:00:50 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	0.00
2	1,4-Dioxane	0.674	0.679	-0.7	97	0.00
3	n-Nitrosodimethylamine	0.317	0.280	11.7	99	-0.01
4 S	2-Fluorophenol	1.080	0.994	8.0	91	-0.01
5 S	Phenol-d6	1.357	1.208	11.0	90	0.00
6	bis(2-Chloroethyl)ether	1.257	1.180	6.1	93	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
8 S	Nitrobenzene-d5	0.250	0.223	10.8	94	0.00
9	Naphthalene	4.126	4.056	1.7	93	0.00
10	Hexachlorobutadiene	0.133	0.133	0.0	94	0.00
11 SURR	2-Methylnaphthalene-d10	0.599	0.578	3.5	93	0.00
12	2-Methylnaphthalene	0.678	0.649	4.3	93	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
14 S	2,4,6-Tribromophenol	0.091	0.080	12.1	93	0.00
15 S	2-Fluorobiphenyl	1.451	1.399	3.6	93	0.00
16	Acenaphthylene	1.699	1.568	7.7	94	0.00
17	Acenaphthene	1.138	1.075	5.5	95	0.00
18	Fluorene	1.454	1.369	5.8	96	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
20	4-Bromophenyl-phenylether	0.198	0.180	9.1	96	0.00
21	Hexachlorobenzene	0.167	0.160	4.2	97	0.00
22	Pentachlorophenol	0.046	0.027	41.3#	83	-0.01
23	Phenanthrene	1.138	1.106	2.8	97	0.00
24	Anthracene	0.979	0.889	9.2	99	0.00
25 SURR	Fluoranthene-d10	4.934	4.546	7.9	99	0.00
26	Fluoranthene	1.401	1.331	5.0	99	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	100	-0.01
28	Pvrene	1.025	1.032	-0.7	100	0.00
29 S	Terphenyl-d14	0.884	0.853	3.5	101	0.00
30	Benzo(a)anthracene	1.118	1.064	4.8	99	0.00
31	Chrysene	1.196	1.176	1.7	99	-0.01
32	Bis(2-ethylhexyl)phthalate	2.136	1.750	18.1	95	0.00
33	Indeno(1,2,3-cd)pyrene	1.409	1.353	4.0	96	0.00
34 I	Perylene-d12	1.000	1.000	0.0	99	0.00
35	Benzo(b)fluoranthene	1.084	0.969	10.6	98	0.00
36	Benzo(k)fluoranthene	1.232	1.153	6.4	98	0.00
37 C	Benzo(a)pyrene	0.995	0.903	9.2	98	0.00
38	Dibenzo(a,h)anthracene	1.047	0.942	10.0	96	0.00
39	Benzo(a,h,i)perylene	1.060	0.981	7.5	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			