

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061719\
 Data File : BE100067.D
 Acq On : 17 Jun 2019 17:25
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICSVBE061719

Quant Time: Jun 17 18:06:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 17 17:23:36 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	88	0.01
2	1,4-Dioxane	0.698	0.654	6.3	83	0.01
3	n-Nitrosodimethylamine	0.289	0.185	36.0#	77	0.02
4 S	2-Fluorophenol	1.083	0.998	7.8	87	0.01
5 S	Phenol-d6	1.306	1.292	1.1	92	0.02
6	bis(2-Chloroethyl)ether	1.225	1.077	12.1	92	0.02
7 I	Naphthalene-d8	1.000	1.000	0.0	92	0.01
8 S	Nitrobenzene-d5	0.396	0.347	12.4	89	0.03
9	Naphthalene	4.423	4.544	-2.7	94	0.01
10	Hexachlorobutadiene	0.380	0.405	-6.6	97	0.03
11 SURR	2-Methylnaphthalene-d10	0.723	0.748	-3.5	98	0.02
12	2-Methylnaphthalene	0.693	0.706	-1.9	97	0.02
13 I	Acenaphthene-d10	1.000	1.000	0.0	96	0.02
14 S	2,4,6-Tribromophenol	0.245	0.220	10.2	99	0.01
15 S	2-Fluorobiphenyl	1.551	1.558	-0.5	95	0.02
16	Acenaphthylene	2.004	1.995	0.4	97	0.02
17	Acenaphthene	1.084	1.083	0.1	95	0.00
18	Fluorene	1.614	1.554	3.7	93	0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	93	0.01
20	4-Bromophenyl-phenylether	0.241	0.236	2.1	93	0.01
21	Hexachlorobenzene	0.247	0.249	-0.8	96	0.01
22	Pentachlorophenol	0.077	0.048	37.7#	69	0.03
23	Phenanthrene	0.975	0.944	3.2	93	0.00
24	Anthracene	0.857	0.790	7.8	91	0.00
25 SURR	Fluoranthene-d10	5.626	5.420	3.7	93	0.00
26	Fluoranthene	1.604	1.565	2.4	93	0.01
27 I	Chrysene-d12	1.000	1.000	0.0	91	0.01
28	Pvrene	0.995	1.117	-12.3	92	0.01
29 S	Terphenyl-d14	0.674	0.723	-7.3	90	0.02
30	Benzo(a)anthracene	1.152	1.285	-11.5	95	0.00
31	Chrysene	1.299	1.417	-9.1	88	0.01
32	Bis(2-ethylhexyl)phthalate	1.604	1.456	9.2	82	0.00
33	Indeno(1,2,3-cd)pyrene	1.535	1.685	-9.8	90	0.01
34 I	Perylene-d12	1.000	1.000	0.0	90	0.00
35	Benzo(b)fluoranthene	1.153	1.152	0.1	89	0.01
36	Benzo(k)fluoranthene	1.249	1.286	-3.0	91	0.00
37 C	Benzo(a)pyrene	1.144	1.173	-2.5	92	0.00
38	Dibenzo(a,h)anthracene	1.116	1.089	2.4	88	0.02
39	Benzo(a,h,i)perylene	1.181	1.203	-1.9	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			