

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061919\
 Data File : BE100085.D
 Acq On : 19 Jun 2019 16:14
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC00.4

Quant Time: Jun 20 14:43:11 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 19 15:07:24 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	92	0.00
2	1,4-Dioxane	0.642	0.724	-12.8	117	0.00
3	n-Nitrosodimethylamine	0.238	0.099	58.4#	130	0.01
4 S	2-Fluorophenol	0.830	0.979	-18.0	115	0.00
5 S	Phenol-d6	1.140	1.214	-6.5	97	0.00
6	bis(2-Chloroethyl)ether	1.119	1.062	5.1	98	-0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	94	0.00
8 S	Nitrobenzene-d5	0.388	0.368	5.2	88	0.01
9	Naphthalene	4.347	4.347	0.0	94	0.00
10	Hexachlorobutadiene	0.427	0.440	-3.0	99	0.00
11 SURR	2-Methylnaphthalene-d10	0.761	0.716	5.9	92	0.00
12	2-Methylnaphthalene	0.715	0.665	7.0	91	0.01
13 I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
14 S	2,4,6-Tribromophenol	0.221	0.200	9.5	98	0.00
15 S	2-Fluorobiphenyl	1.573	1.607	-2.2	91	0.00
16	Acenaphthylene	1.899	1.819	4.2	92	0.00
17	Acenaphthene	1.062	1.045	1.6	93	0.01
18	Fluorene	1.581	1.546	2.2	93	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	92	0.00
20	4-Bromophenyl-phenylether	0.238	0.238	0.0	96	0.00
21	Hexachlorobenzene	0.262	0.269	-2.7	96	0.00
22	Pentachlorophenol	0.053	0.038	28.3#	92	0.00
23	Phenanthrene	0.958	0.938	2.1	92	0.00
24	Anthracene	0.808	0.730	9.7	90	0.01
25 SURR	Fluoranthene-d10	5.718	5.686	0.6	93	0.00
26	Fluoranthene	1.538	1.522	1.0	93	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	95	0.01
28	Pvrene	1.028	0.995	3.2	93	0.00
29 S	Terphenyl-d14	0.800	0.780	2.5	93	0.00
30	Benzo(a)anthracene	1.190	1.145	3.8	100	0.00
31	Chrysene	1.383	1.362	1.5	91	0.00
32	Bis(2-ethylhexyl)phthalate	1.123	0.936	16.7	92	0.01
33	Indeno(1,2,3-cd)pyrene	1.556	1.518	2.4	90	0.00
34 I	Perylene-d12	1.000	1.000	0.0	94	0.00
35	Benzo(b)fluoranthene	1.174	1.126	4.1	91	0.00
36	Benzo(k)fluoranthene	1.335	1.346	-0.8	98	0.00
37 C	Benzo(a)pyrene	1.133	1.095	3.4	91	0.00
38	Dibenzo(a,h)anthracene	1.161	1.091	6.0	91	0.00
39	Benzo(a,h,i)perylene	1.207	1.181	2.2	92	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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