

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031419\
 Data File : BE098965.D
 Acq On : 15 Mar 2019 4:31
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC00.4

Quant Time: Mar 15 06:06:51 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 15 05:31:33 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	120	0.00
2	1,4-Dioxane	0.789	0.705	10.6	111	0.00
3	n-Nitrosodimethylamine	1.006	0.783	22.2	92	0.01
4 S	2-Fluorophenol	1.196	1.095	8.4	106	0.01
5 S	Phenol-d6	1.689	1.633	3.3	116	0.00
6	bis(2-Chloroethyl)ether	1.304	1.158	11.2	111	0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	130	0.00
8 S	Nitrobenzene-d5	0.419	0.335	20.0	115	0.00
9	Naphthalene	4.581	4.484	2.1	130	-0.01
10	Hexachlorobutadiene	0.303	0.287	5.3	127	0.00
11 SURR	2-Methylnaphthalene-d10	0.739	0.713	3.5	128	0.00
12	2-Methylnaphthalene	0.743	0.723	2.7	130	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	135	0.00
14 S	2,4,6-Tribromophenol	0.202	0.158	21.8	107	0.00
15 S	2-Fluorobiphenyl	1.656	1.567	5.4	129	-0.01
16	Acenaphthylene	2.057	1.971	4.2	136	-0.01
17	Acenaphthene	1.220	1.194	2.1	133	-0.01
18	Fluorene	1.760	1.625	7.7	128	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	120	0.00
20	4-Bromophenyl-phenylether	0.217	0.208	4.1	117	0.00
21	Hexachlorobenzene	0.258	0.247	4.3	121	0.00
22	Pentachlorophenol	0.052	0.041	21.2	127	0.00
23	Phenanthrene	1.137	1.059	6.9	114	0.00
24	Anthracene	0.972	0.833	14.3	109	0.00
25 SURR	Fluoranthene-d10	6.385	5.384	15.7	105	0.00
26	Fluoranthene	1.605	1.356	15.5	105	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	102	-0.01
28	Pvrene	1.224	1.246	-1.8	104	0.00
29 S	Terphenyl-d14	0.972	0.926	4.7	98	0.00
30	Benzo(a)anthracene	1.248	1.090	12.7	91	0.00
31	Chrysene	1.336	1.336	0.0	105	-0.01
32	Bis(2-ethylhexyl)phthalate	2.657	2.306	13.2	97	-0.01
33	Indeno(1,2,3-cd)pyrene	1.410	1.321	6.3	97	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	106	-0.01
35	Benzo(b)fluoranthene	1.315	1.255	4.6	105	0.00
36	Benzo(k)fluoranthene	1.377	1.311	4.8	106	-0.01
37 C	Benzo(a)pyrene	1.261	1.235	2.1	108	-0.01
38	Dibenzo(a,h)anthracene	1.193	0.930	22.0	84	-0.01
39	Benzo(a,h,i)perylene	1.248	1.145	8.3	100	-0.01

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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