

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031319\
 Data File : BE098909.D
 Acq On : 13 Mar 2019 13:28
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC00.4

Quant Time: Mar 14 08:56:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 06 15:33:24 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	95	0.00
2	1,4-Dioxane	0.789	0.799	-1.3	100	0.00
3	n-Nitrosodimethylamine	1.006	0.822	18.3	77	0.04
4 S	2-Fluorophenol	1.196	0.888	25.8#	68	0.01
5 S	Phenol-d6	1.689	1.564	7.4	88	0.01
6	bis(2-Chloroethyl)ether	1.304	0.947	27.4#	72	0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
8 S	Nitrobenzene-d5	0.419	0.342	18.4	92	0.01
9	Naphthalene	4.581	4.439	3.1	100	0.00
10	Hexachlorobutadiene	0.303	0.300	1.0	103	0.00
11 SURR	2-Methylnaphthalene-d10	0.739	0.692	6.4	96	0.00
12	2-Methylnaphthalene	0.743	0.673	9.4	94	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
14 S	2,4,6-Tribromophenol	0.202	0.152	24.8	81	0.00
15 S	2-Fluorobiphenyl	1.656	1.552	6.3	101	0.00
16	Acenaphthylene	2.057	1.928	6.3	105	0.00
17	Acenaphthene	1.220	1.158	5.1	102	-0.02
18	Fluorene	1.760	1.618	8.1	101	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	97	0.00
20	4-Bromophenyl-phenylether	0.217	0.207	4.6	95	0.00
21	Hexachlorobenzene	0.258	0.252	2.3	100	0.00
22	Pentachlorophenol	0.052	0.030	42.3#	77	0.01
23	Phenanthrene	1.137	1.046	8.0	92	0.00
24	Anthracene	0.972	0.791	18.6	84	0.00
25 SURR	Fluoranthene-d10	6.385	5.582	12.6	88	0.00
26	Fluoranthene	1.605	1.392	13.3	88	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	91	-0.01
28	Pvrene	1.224	1.176	3.9	88	0.00
29 S	Terphenyl-d14	0.972	0.878	9.7	83	0.00
30	Benzo(a)anthracene	1.248	1.020	18.3	76	0.00
31	Chrysene	1.336	1.312	1.8	91	-0.01
32	Bis(2-ethylhexyl)phthalate	2.657	1.896	28.6#	71	-0.01
33	Indeno(1,2,3-cd)pyrene	1.410	1.331	5.6	87	0.00
34 I	Perylene-d12	1.000	1.000	0.0	95	-0.01
35	Benzo(b)fluoranthene	1.315	1.185	9.9	89	0.00
36	Benzo(k)fluoranthene	1.377	1.264	8.2	92	-0.01
37 C	Benzo(a)pyrene	1.261	1.171	7.1	91	-0.01
38	Dibenzo(a,h)anthracene	1.193	1.035	13.2	84	-0.01
39	Benzo(a,h,i)perylene	1.248	1.240	0.6	97	-0.81#

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