

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031419\
 Data File : BE098946.D
 Acq On : 14 Mar 2019 15:41
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC00.4

Quant Time: Mar 15 05:45:45 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	99	0.01
2	1,4-Dioxane	0.789	0.775	1.8	101	0.00
3	n-Nitrosodimethylamine	1.006	0.887	11.8	85	0.04
4 S	2-Fluorophenol	1.196	0.817	31.7#	65	0.03
5 S	Phenol-d6	1.689	1.384	18.1	81	0.03
6	bis(2-Chloroethyl)ether	1.304	0.980	24.8	77	0.04
7 I	Naphthalene-d8	1.000	1.000	0.0	112	0.01
8 S	Nitrobenzene-d5	0.419	0.316	24.6	94	0.03
9	Naphthalene	4.581	4.478	2.2	113	0.01
10	Hexachlorobutadiene	0.303	0.290	4.3	111	0.01
11 SURR	2-Methylnaphthalene-d10	0.739	0.729	1.4	113	0.02
12	2-Methylnaphthalene	0.743	0.660	11.2	103	0.02
13 I	Acenaphthene-d10	1.000	1.000	0.0	117	0.02
14 S	2,4,6-Tribromophenol	0.202	0.129	36.1#	76	0.01
15 S	2-Fluorobiphenyl	1.656	1.539	7.1	110	0.02
16	Acenaphthylene	2.057	1.861	9.5	111	0.00
17	Acenaphthene	1.220	1.146	6.1	110	0.00
18	Fluorene	1.760	1.541	12.4	105	0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
20	4-Bromophenyl-phenylether	0.217	0.210	3.2	100	0.01
21	Hexachlorobenzene	0.258	0.247	4.3	103	0.00
22	Pentachlorophenol	0.052	0.022	57.7#	57	0.01
23	Phenanthrene	1.137	1.059	6.9	97	0.00
24	Anthracene	0.972	0.765	21.3	85	0.01
25 SURR	Fluoranthene-d10	6.385	5.411	15.3	89	0.00
26	Fluoranthene	1.605	1.347	16.1	89	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	83	0.00
28	Pvrene	1.224	1.289	-5.3	88	0.00
29 S	Terphenyl-d14	0.972	0.908	6.6	78	0.01
30	Benzo(a)anthracene	1.248	1.031	17.4	70	0.01
31	Chrysene	1.336	1.316	1.5	84	0.00
32	Bis(2-ethylhexyl)phthalate	2.657	1.836	30.9#	63	0.00
33	Indeno(1,2,3-cd)pyrene	1.410	1.130	19.9	68	0.01
34 I	Perylene-d12	1.000	1.000	0.0	84	0.00
35	Benzo(b)fluoranthene	1.315	1.186	9.8	78	0.00
36	Benzo(k)fluoranthene	1.377	1.372	0.4	88	0.00
37 C	Benzo(a)pyrene	1.261	1.161	7.9	80	0.00
38	Dibenzo(a,h)anthracene	1.193	0.898	24.7	64	0.01
39	Benzo(a,h,i)perylene	1.248	1.115	10.7	76	0.01

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