

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031219\
 Data File : BE098890.D
 Acq On : 12 Mar 2019 10:14
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC00.4

Quant Time: Mar 12 12:30:54 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	116	0.00
2	1,4-Dioxane	0.789	0.778	1.4	119	0.00
3	n-Nitrosodimethylamine	1.006	0.454	54.9#	52	0.04
4 S	2-Fluorophenol	1.196	1.043	12.8	98	0.01
5 S	Phenol-d6	1.689	1.651	2.2	114	0.00
6	bis(2-Chloroethyl)ether	1.304	0.988	24.2	91	0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	126	0.00
8 S	Nitrobenzene-d5	0.419	0.336	19.8	112	0.00
9	Naphthalene	4.581	4.497	1.8	127	0.00
10	Hexachlorobutadiene	0.303	0.287	5.3	123	0.00
11 SURR	2-Methylnaphthalene-d10	0.739	0.715	3.2	125	0.00
12	2-Methylnaphthalene	0.743	0.686	7.7	120	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	129	0.00
14 S	2,4,6-Tribromophenol	0.202	0.172	14.9	112	0.00
15 S	2-Fluorobiphenyl	1.656	1.649	0.4	130	0.00
16	Acenaphthylene	2.057	1.965	4.5	130	0.00
17	Acenaphthene	1.220	1.207	1.1	129	-0.01
18	Fluorene	1.760	1.674	4.9	126	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	121	0.00
20	4-Bromophenyl-phenylether	0.217	0.199	8.3	112	0.00
21	Hexachlorobenzene	0.258	0.251	2.7	124	0.00
22	Pentachlorophenol	0.052	0.031	40.4#	98	0.01
23	Phenanthrene	1.137	1.080	5.0	117	0.00
24	Anthracene	0.972	0.870	10.5	114	0.00
25 SURR	Fluoranthene-d10	6.385	5.673	11.2	111	0.00
26	Fluoranthene	1.605	1.411	12.1	110	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	108	-0.01
28	Pvrene	1.224	1.245	-1.7	111	0.00
29 S	Terphenyl-d14	0.972	0.939	3.4	105	0.00
30	Benzo(a)anthracene	1.248	1.095	12.3	97	0.00
31	Chrysene	1.336	1.350	-1.0	112	0.00
32	Bis(2-ethylhexyl)phthalate	2.657	2.184	17.8	97	-0.01
33	Indeno(1,2,3-cd)pyrene	1.410	1.411	-0.1	110	0.00
34 I	Perylene-d12	1.000	1.000	0.0	112	-0.01
35	Benzo(b)fluoranthene	1.315	1.294	1.6	114	0.00
36	Benzo(k)fluoranthene	1.377	1.356	1.5	116	-0.01
37 C	Benzo(a)pyrene	1.261	1.227	2.7	113	0.00
38	Dibenzo(a,h)anthracene	1.193	1.128	5.4	107	-0.02
39	Benzo(a,h,i)perylene	1.248	1.205	3.4	110	-0.01

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031219\
Data File : BE098890.D
Acq On : 12 Mar 2019 10:14
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Mar 12 12:30:54 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	AvgRF	CCRF	%Dev Area	% Dev(min)

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