

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121118\
 Data File : BE098360.D
 Acq On : 12 Dec 2018 13:38
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 12 14:10:39 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 10 14:59:55 2018
 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	135	0.00
2	1,4-Dioxane	0.692	0.689	0.4	137	0.00
3	n-Nitrosodimethylamine	0.623	0.554	11.1	129	0.00
4 S	2-Fluorophenol	0.936	0.960	-2.6	142	0.00
5 S	Phenol-d6	1.326	1.345	-1.4	136	0.00
6	bis(2-Chloroethyl)ether	1.260	1.123	10.9	124	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	128	0.01
8 S	Nitrobenzene-d5	0.364	0.405	-11.3	155#	0.03
9	Naphthalene	4.025	3.926	2.5	128	0.01
10	Hexachlorobutadiene	0.256	0.258	-0.8	132	0.01
11 SURR	2-Methylnaphthalene-d10	0.650	0.611	6.0	121	0.01
12	2-Methylnaphthalene	0.654	0.612	6.4	122	0.01
13 I	Acenaphthene-d10	1.000	1.000	0.0	117	0.00
14 S	2,4,6-Tribromophenol	0.147	0.145	1.4	122	0.01
15 S	2-Fluorobiphenyl	1.687	1.640	2.8	125	0.01
16	Acenaphthylene	1.874	1.804	3.7	117	0.01
17	Acenaphthene	1.126	1.093	2.9	115	0.01
18	Fluorene	1.506	1.420	5.7	111	0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
20	4-Bromophenyl-phenylether	0.215	0.206	4.2	110	0.01
21	Hexachlorobenzene	0.231	0.229	0.9	111	0.00
22	Pentachlorophenol	0.046	0.045	2.2	135	0.00
23	Phenanthrene	0.972	0.946	2.7	108	0.00
24	Anthracene	0.866	0.800	7.6	103	0.01
25 SURR	Fluoranthene-d10	5.131	4.997	2.6	103	0.00
26	Fluoranthene	1.286	1.255	2.4	103	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	127	0.00
28	Pvrene	1.255	1.089	13.2	108	0.00
29 S	Terphenyl-d14	0.972	0.814	16.3	108	0.00
30	Benzo(a)anthracene	1.139	1.032	9.4	118	0.00
31	Chrysene	1.193	1.153	3.4	124	0.00
32	Bis(2-ethylhexyl)phthalate	2.313	2.149	7.1	119	0.00
33	Indeno(1,2,3-cd)pyrene	1.187	1.166	1.8	132	0.00
34 I	Perylene-d12	1.000	1.000	0.0	134	0.00
35	Benzo(b)fluoranthene	1.094	1.042	4.8	132	0.00
36	Benzo(k)fluoranthene	1.274	1.280	-0.5	140	0.00
37 C	Benzo(a)pyrene	1.129	1.093	3.2	132	0.00
38	Dibenzo(a,h)anthracene	1.063	1.029	3.2	133	0.00
39	Benzo(a,h,i)perylene	1.071	1.020	4.8	131	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121118\
Data File : BE098360.D
Acq On : 12 Dec 2018 13:38
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Dec 12 14:10:39 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 10 14:59:55 2018
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			