

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121318\
 Data File : BE098416.D
 Acq On : 14 Dec 2018 2:58
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 14 10:06:54 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	72	0.00
2	1,4-Dioxane	0.692	0.677	2.2	72	-0.01
3	n-Nitrosodimethylamine	0.623	0.350	43.8#	44#	0.01
4 S	2-Fluorophenol	0.936	0.907	3.1	72	0.00
5 S	Phenol-d6	1.326	1.348	-1.7	73	0.00
6	bis(2-Chloroethyl)ether	1.260	1.153	8.5	68	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	74	-0.01
8 S	Nitrobenzene-d5	0.364	0.317	12.9	71	0.01
9	Naphthalene	4.025	3.861	4.1	73	0.00
10	Hexachlorobutadiene	0.256	0.260	-1.6	77	0.00
11 SURR	2-Methylnaphthalene-d10	0.650	0.605	6.9	70	0.00
12	2-Methylnaphthalene	0.654	0.596	8.9	69	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	72	0.00
14 S	2,4,6-Tribromophenol	0.147	0.136	7.5	70	0.00
15 S	2-Fluorobiphenyl	1.687	1.536	9.0	72	0.00
16	Acenaphthylene	1.874	1.807	3.6	72	0.00
17	Acenaphthene	1.126	1.064	5.5	69	0.00
18	Fluorene	1.506	1.437	4.6	69	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	71	-0.01
20	4-Bromophenyl-phenylether	0.215	0.187	13.0	64	0.00
21	Hexachlorobenzene	0.231	0.220	4.8	68	0.00
22	Pentachlorophenol	0.046	0.018	60.9#	36#	0.01
23	Phenanthrene	0.972	0.926	4.7	68	0.00
24	Anthracene	0.866	0.745	14.0	61	0.00
25 SURR	Fluoranthene-d10	5.131	5.378	-4.8	71	0.00
26	Fluoranthene	1.286	1.374	-6.8	72	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	86	0.00
28	Pvrene	1.255	1.072	14.6	72	0.00
29 S	Terphenyl-d14	0.972	0.792	18.5	71	0.00
30	Benzo(a)anthracene	1.139	1.024	10.1	79	-0.01
31	Chrysene	1.193	1.116	6.5	81	0.00
32	Bis(2-ethylhexyl)phthalate	2.313	1.851	20.0	69	-0.01
33	Indeno(1,2,3-cd)pyrene	1.187	1.120	5.6	85	0.00
34 I	Perylene-d12	1.000	1.000	0.0	88	-0.01
35	Benzo(b)fluoranthene	1.094	1.085	0.8	90	0.00
36	Benzo(k)fluoranthene	1.274	1.313	-3.1	94	-0.01
37 C	Benzo(a)pyrene	1.129	1.119	0.9	89	-0.01
38	Dibenzo(a,h)anthracene	1.063	0.979	7.9	83	-0.01
39	Benzo(a,h,i)perylene	1.071	1.026	4.2	87	-0.02

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121318\
Data File : BE098416.D
Acq On : 14 Dec 2018 2:58
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Dec 14 10:06:54 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			