

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121318\
 Data File : BE098397.D
 Acq On : 13 Dec 2018 15:16
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 13 20:57:25 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	80	0.00
2	1,4-Dioxane	0.692	0.635	8.2	75	0.00
3	n-Nitrosodimethylamine	0.623	0.377	39.5#	52	0.01
4 S	2-Fluorophenol	0.936	0.887	5.2	78	0.01
5 S	Phenol-d6	1.326	1.298	2.1	78	0.00
6	bis(2-Chloroethyl)ether	1.260	1.016	19.4	67	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	79	-0.01
8 S	Nitrobenzene-d5	0.364	0.347	4.7	82	0.01
9	Naphthalene	4.025	3.918	2.7	79	0.00
10	Hexachlorobutadiene	0.256	0.260	-1.6	82	0.00
11 SURR	2-Methylnaphthalene-d10	0.650	0.608	6.5	75	0.00
12	2-Methylnaphthalene	0.654	0.587	10.2	72	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	79	0.00
14 S	2,4,6-Tribromophenol	0.147	0.142	3.4	80	0.00
15 S	2-Fluorobiphenyl	1.687	1.532	9.2	79	0.00
16	Acenaphthylene	1.874	1.796	4.2	78	0.00
17	Acenaphthene	1.126	1.066	5.3	75	0.00
18	Fluorene	1.506	1.442	4.2	76	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	76	-0.01
20	4-Bromophenyl-phenylether	0.215	0.184	14.4	68	0.00
21	Hexachlorobenzene	0.231	0.221	4.3	74	0.00
22	Pentachlorophenol	0.046	0.029	37.0#	60	0.00
23	Phenanthrene	0.972	0.919	5.5	73	0.00
24	Anthracene	0.866	0.737	14.9	66	0.00
25 SURR	Fluoranthene-d10	5.131	5.074	1.1	72	0.00
26	Fluoranthene	1.286	1.294	-0.6	73	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	88	0.00
28	Pvrene	1.255	1.082	13.8	74	0.00
29 S	Terphenyl-d14	0.972	0.768	21.0	70	0.00
30	Benzo(a)anthracene	1.139	1.029	9.7	81	-0.01
31	Chrysene	1.193	1.110	7.0	82	0.00
32	Bis(2-ethylhexyl)phthalate	2.313	1.748	24.4	67	-0.01
33	Indeno(1,2,3-cd)pyrene	1.187	1.145	3.5	89	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	93	-0.01
35	Benzo(b)fluoranthene	1.094	1.039	5.0	91	0.00
36	Benzo(k)fluoranthene	1.274	1.289	-1.2	98	-0.01
37 C	Benzo(a)pyrene	1.129	1.059	6.2	89	-0.01
38	Dibenzo(a,h)anthracene	1.063	0.981	7.7	88	-0.01
39	Benzo(a,h,i)perylene	1.071	0.999	6.7	89	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121318\
Data File : BE098397.D
Acq On : 13 Dec 2018 15:16
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Dec 13 20:57:25 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			