

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE071918\
 Data File : BE097014.D
 Acq On : 19 Jul 2018 20:48
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 20 04:33:44 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 17 15:02:56 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	67	0.00
2	1,4-Dioxane	0.573	0.580	-1.2	70	0.00
3	n-Nitrosodimethylamine	0.644	0.698	-8.4	77	0.00
4 S	2-Fluorophenol	0.957	1.020	-6.6	81	0.00
5 S	Phenol-d6	1.460	1.464	-0.3	73	0.00
6	bis(2-Chloroethyl)ether	1.433	1.374	4.1	67	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	61	0.00
8 S	Nitrobenzene-d5	0.292	0.311	-6.5	70	0.00
9	Naphthalene	3.581	3.366	6.0	60	0.00
10	Hexachlorobutadiene	0.145	0.154	-6.2	68	0.01
11 SURR	2-Methylnaphthalene-d10	0.583	0.542	7.0	60	0.00
12	2-Methylnaphthalene	0.607	0.572	5.8	60	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	70	0.00
14 S	2,4,6-Tribromophenol	0.101	0.109	-7.9	93	0.00
15 S	2-Fluorobiphenyl	1.488	1.309	12.0	63	0.00
16	Acenaphthylene	1.827	1.676	8.3	67	0.01
17	Acenaphthene	1.136	1.006	11.4	65	0.00
18	Fluorene	1.275	1.298	-1.8	76	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	85	0.00
20	4-Bromophenyl-phenylether	0.189	0.179	5.3	85	0.00
21	Hexachlorobenzene	0.210	0.197	6.2	83	0.00
22	Pentachlorophenol	0.034	0.044	-29.4#	123	0.00
23	Phenanthrene	0.958	0.899	6.2	84	0.00
24	Anthracene	0.834	0.785	5.9	86	0.00
25 SURR	Fluoranthene-d10	3.480	3.760	-8.0	100	0.00
26	Fluoranthene	1.010	1.034	-2.4	93	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	106	0.00
28	Pvrene	1.181	1.035	12.4	96	0.00
29 S	Terphenyl-d14	0.772	0.715	7.4	103	0.00
30	Benzo(a)anthracene	0.952	0.924	2.9	112	0.00
31	Chrysene	1.068	0.998	6.6	101	0.00
32	Bis(2-ethylhexyl)phthalate	0.951	0.939	1.3	120	0.00
33	Indeno(1,2,3-cd)pyrene	0.778	0.977	-25.6#	149	0.00
34 I	Perylene-d12	1.000	1.000	0.0	121	0.00
35	Benzo(b)fluoranthene	1.021	0.877	14.1	112	0.00
36	Benzo(k)fluoranthene	1.214	1.016	16.3	105	0.00
37 C	Benzo(a)pyrene	0.913	0.772	15.4	113	0.00
38	Dibenzo(a,h)anthracene	0.746	0.882	-18.2	162#	0.00
39	Benzo(a,h,i)perylene	0.830	0.904	-8.9	145	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE071918\
Data File : BE097014.D
Acq On : 19 Jul 2018 20:48
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Jul 20 04:33:44 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 17 15:02:56 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		