

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE081618\
 Data File : BE097286.D
 Acq On : 16 Aug 2018 11:21
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 16 15:03:49 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE081518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 15 14:23:39 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)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	94	0.00
2	1,4-Dioxane	0.449	0.400	10.9	87	0.00
3	n-Nitrosodimethylamine	0.521	0.435	16.5	96	0.00
4 S	2-Fluorophenol	1.315	1.168	11.2	90	-0.01
5 S	Phenol-d6	1.879	1.709	9.0	90	0.00
6	bis(2-Chloroethyl)ether	1.497	1.362	9.0	91	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
8 S	Nitrobenzene-d5	0.366	0.331	9.6	90	0.00
9	Naphthalene	3.740	3.441	8.0	91	-0.01
10	Hexachlorobutadiene	0.169	0.164	3.0	96	0.00
11 SURR	2-Methylnaphthalene-d10	0.581	0.532	8.4	90	0.00
12	2-Methylnaphthalene	0.612	0.559	8.7	90	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	92	0.00
14 S	2,4,6-Tribromophenol	0.134	0.106	20.9	82	0.00
15 S	2-Fluorobiphenyl	1.503	1.444	3.9	92	0.00
16	Acenaphthylene	1.846	1.763	4.5	90	0.00
17	Acenaphthene	1.061	1.025	3.4	90	0.00
18	Fluorene	1.317	1.235	6.2	90	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	90	0.00
20	4-Bromophenyl-phenylether	0.209	0.202	3.3	90	0.00
21	Hexachlorobenzene	0.237	0.239	-0.8	92	0.00
22	Pentachlorophenol	0.053	0.041	22.6	93	0.00
23	Phenanthrene	0.965	0.931	3.5	90	0.00
24	Anthracene	0.865	0.776	10.3	86	0.00
25 SURR	Fluoranthene-d10	3.998	3.636	9.1	90	0.00
26	Fluoranthene	1.068	0.977	8.5	89	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	86	0.00
28	Pyrene	1.079	1.044	3.2	85	0.00
29 S	Terphenyl-d14	0.752	0.726	3.5	85	0.00
30	Benzo(a)anthracene	1.006	0.925	8.1	81	-0.01
31	Chrysene	1.030	1.019	1.1	88	0.00
32	Bis(2-ethylhexyl)phthalate	1.301	0.856	34.2#	61	0.00
33	Indeno(1,2,3-cd)pyrene	1.109	1.017	8.3	80	0.00
34 I	Perylene-d12	1.000	1.000	0.0	82	0.00
35	Benzo(b)fluoranthene	0.995	0.930	6.5	80	0.00
36	Benzo(k)fluoranthene	1.210	1.255	-3.7	87	0.00
37 C	Benzo(a)pyrene	0.999	0.945	5.4	80	0.00
38	Dibenzo(a,h)anthracene	0.986	0.930	5.7	82	-0.01
39	Benzo(g,h,i)perylene	0.995	0.903	9.2	78	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE081618\
Data File : BE097286.D
Acq On : 16 Aug 2018 11:21
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Aug 16 15:03:49 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE081518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 15 14:23:39 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			