

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE071918\
 Data File : BE097000.D
 Acq On : 19 Jul 2018 12:04
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 19 13:31:13 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	63	0.00
2	1,4-Dioxane	0.573	0.613	-7.0	70	0.00
3	n-Nitrosodimethylamine	0.644	0.669	-3.9	69	0.00
4 S	2-Fluorophenol	0.957	0.970	-1.4	72	0.00
5 S	Phenol-d6	1.460	1.368	6.3	64	0.00
6	bis(2-Chloroethyl)ether	1.433	1.387	3.2	63	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	55	0.00
8 S	Nitrobenzene-d5	0.292	0.309	-5.8	63	0.00
9	Naphthalene	3.581	3.442	3.9	56	0.00
10	Hexachlorobutadiene	0.145	0.155	-6.9	62	0.01
11 SURR	2-Methylnaphthalene-d10	0.583	0.545	6.5	55	0.00
12	2-Methylnaphthalene	0.607	0.579	4.6	55	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	65	0.00
14 S	2,4,6-Tribromophenol	0.101	0.103	-2.0	81	0.00
15 S	2-Fluorobiphenyl	1.488	1.321	11.2	59	0.01
16	Acenaphthylene	1.827	1.655	9.4	61	0.01
17	Acenaphthene	1.136	0.998	12.1	60	0.00
18	Fluorene	1.275	1.284	-0.7	69	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	79	0.01
20	4-Bromophenyl-phenylether	0.189	0.176	6.9	78	0.01
21	Hexachlorobenzene	0.210	0.193	8.1	75	0.00
22	Pentachlorophenol	0.034	0.047	-38.2#	122	0.01
23	Phenanthrene	0.958	0.913	4.7	79	0.00
24	Anthracene	0.834	0.778	6.7	79	0.00
25 SURR	Fluoranthene-d10	3.480	3.699	-6.3	91	0.00
26	Fluoranthene	1.010	1.013	-0.3	84	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	99	0.00
28	Pvrene	1.181	1.006	14.8	88	0.00
29 S	Terphenyl-d14	0.772	0.711	7.9	96	0.00
30	Benzo(a)anthracene	0.952	0.918	3.6	105	0.00
31	Chrysene	1.068	0.973	8.9	93	0.00
32	Bis(2-ethylhexyl)phthalate	0.951	0.885	6.9	106	0.00
33	Indeno(1,2,3-cd)pyrene	0.778	0.968	-24.4	139	0.00
34 I	Perylene-d12	1.000	1.000	0.0	108	0.00
35	Benzo(b)fluoranthene	1.021	0.914	10.5	104	0.00
36	Benzo(k)fluoranthene	1.214	1.044	14.0	97	0.00
37 C	Benzo(a)pyrene	0.913	0.837	8.3	110	0.00
38	Dibenzo(a,h)anthracene	0.746	0.898	-20.4	148	0.00
39	Benzo(a,h,i)perylene	0.830	0.941	-13.4	136	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

(#) = Out of Range	SPCC's out = 0 CCC's out = 0			