

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM082916\
 Data File : BM007340.D
 Acq On : 30 Aug 2016 11:14
 Operator : UM/SJ
 Sample : SSTDC0020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02037

Quant Time: Aug 30 11:57:48 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM082316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 30 04:03:42 2016
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	87	0.00
2	1,4-Dioxane	0.436	0.444	-1.8	91	0.00
3 S	1,4-Dioxane-d8	0.391	0.420	-7.4	95	0.00
4	Benzaldehyde	1.076	1.146	-6.5	81	0.00
5 S	Phenol-d5	1.889	1.838	2.7	85	0.00
6	Phenol	1.960	1.897	3.2	84	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	0.993	0.965	2.8	82	0.00
8	Bis(2-Chloroethyl)ether	1.365	1.362	0.2	84	0.00
9 S	2-Chlorophenol-d4	1.396	1.405	-0.6	88	0.00
10	2-Chlorophenol	1.430	1.434	-0.3	85	0.00
11	2-Methylphenol	1.532	1.491	2.7	85	0.00
12	2,2'-oxybis(1-Chloropropane	1.536	1.495	2.7	81	0.00
13 S	4-Methylphenol-d8	1.648	1.589	3.6	84	0.00
14	Acetophenone	2.527	2.518	0.4	84	0.00
15 P	N-Nitroso-di-n-propylamine	1.343	1.306	2.8	82	0.00
16	4-Methylphenol	1.718	1.682	2.1	84	0.00
17	Hexachloroethane	0.556	0.579	-4.1	89	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	84	0.00
19 S	Nitrobenzene-d5	0.147	0.152	-3.4	86	0.00
20	Nitrobenzene	0.377	0.394	-4.5	86	0.00
21	Isophorone	0.768	0.763	0.7	80	0.00
22 S	2-Nitrophenol-d4	0.171	0.175	-2.3	84	0.00
23 C	2-Nitrophenol	0.186	0.186	0.0	80	0.00
24	2,4-Dimethylphenol	0.411	0.430	-4.6	85	0.00
25	Bis(2-Chloroethoxy)methane	0.433	0.432	0.2	81	0.00
26 S	2,4-Dichlorophenol-d3	0.341	0.358	-5.0	86	0.00
27 C	2,4-Dichlorophenol	0.339	0.356	-5.0	85	0.00
28	Naphthalene	0.995	1.028	-3.3	84	0.00
29 S	4-Chloroaniline-d4	0.426	0.456	-7.0	84	0.00
30	4-Chloroaniline	0.438	0.467	-6.6	82	0.00
31 C	Hexachlorobutadiene	0.229	0.249	-8.7	89	0.00
32	Caprolactam	0.137	0.134	2.2	77	0.00
33 C	4-Chloro-3-methylphenol	0.420	0.436	-3.8	83	0.00
34	2-Methylnaphthalene	0.814	0.847	-4.1	85	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	83	0.00
36	1,2,4,5-Tetrachlorobenzene	0.666	0.717	-7.7	86	0.00
37	Hexachlorocyclopentadiene	0.402	0.361	10.2	74	0.00
38 C	2,4,6-Trichlorophenol	0.471	0.491	-4.2	83	0.00
39	2,4,5-Trichlorophenol	0.491	0.525	-6.9	86	0.00
40	1,1'-Biphenyl	1.551	1.612	-3.9	83	0.00
41	2-Chloronaphthalene	1.216	1.282	-5.4	84	0.00
42	2-Nitroaniline	0.397	0.424	-6.8	84	0.00
43 S	Dimethylphthalate-d6	1.725	1.738	-0.8	81	0.00
44	Dimethylphthalate	1.719	1.723	-0.2	79	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM082916\
 Data File : BM007340.D
 Acq On : 30 Aug 2016 11:14
 Operator : UM/SJ
 Sample : SSTDC0020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02037

Quant Time: Aug 30 11:57:48 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM082316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 30 04:03:42 2016
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	2,6-Dinitrotoluene	0.351	0.363	-3.4	81	0.00
46 S	Acenaphthylene-d8	2.000	2.099	-5.0	84	0.00
47	Acenaphthylene	2.058	2.154	-4.7	83	0.00
48	3-Nitroaniline	0.359	0.387	-7.8	81	0.00
49 C	Acenaphthene	1.339	1.389	-3.7	83	0.00
50	2,4-Dinitrophenol	0.238	0.188	21.0#	69	0.00
51 S	4-Nitrophenol-d4	0.326	0.330	-1.2	80	0.00
52	4-Nitrophenol	0.347	0.364	-4.9	84	0.00
53	Dibenzofuran	2.004	2.121	-5.8	85	0.00
54	2,4-Dinitrotoluene	0.538	0.561	-4.3	83	0.00
55	2,3,4,6-Tetrachlorophenol	0.504	0.536	-6.3	84	0.00
56	Diethylphthalate	1.793	1.806	-0.7	80	0.00
57 S	Fluorene-d10	1.492	1.563	-4.8	83	0.00
58	Fluorene	1.650	1.744	-5.7	84	0.00
59	4-Chlorophenyl-phenylether	0.863	0.909	-5.3	84	0.00
60	4-Nitroaniline	0.407	0.438	-7.6	82	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	84	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.126	0.111	11.9	73	0.00
63	4,6-Dinitro-2-methylphenol	0.134	0.119	11.2	73	0.00
64	N-Nitrosodiphenylamine	0.558	0.575	-3.0	82	0.00
65	4-Bromophenyl-phenylether	0.226	0.235	-4.0	83	0.00
66	Hexachlorobenzene	0.253	0.263	-4.0	84	0.00
67	Atrazine	0.236	0.244	-3.4	79	0.00
68 C	Pentachlorophenol	0.163	0.161	1.2	81	0.00
69	Phenanthrene	1.074	1.117	-4.0	84	0.00
70 S	Anthracene-d10	0.934	0.974	-4.3	84	0.00
71	Anthracene	1.108	1.151	-3.9	83	0.00
72	Carbazole	0.963	1.027	-6.6	82	0.00
73	Di-n-butylphthalate	1.212	1.207	0.4	79	0.00
74 C	Fluoranthene	1.347	1.474	-9.4	83	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	81	0.00
76 S	Pyrene-d10	0.836	0.883	-5.6	83	0.00
77	Pyrene	1.073	1.137	-6.0	82	0.00
78	Butylbenzylphthalate	0.446	0.457	-2.5	80	0.00
79	3,3'-Dichlorobenzidine	0.397	0.406	-2.3	72	0.00
80	Benzo(a)anthracene	1.179	1.223	-3.7	81	0.00
81	Bis(2-ethylhexyl)phthalate	0.648	0.662	-2.2	79	0.00
82	Chrysene	1.068	1.104	-3.4	80	0.00
83 I	Perylene-d12	1.000	1.000	0.0	74	0.00
84	Di-n-octyl phthalate	0.996	1.176	-18.1	78	0.00
85	Benzo(b)fluoranthene	1.183	1.275	-7.8	75	0.00
86	Benzo(k)fluoranthene	1.093	1.158	-5.9	76	0.00
87 S	Benzo(a)pyrene-d12	0.921	0.946	-2.7	74	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM082916\
Data File : BM007340.D
Acq On : 30 Aug 2016 11:14
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02037

Quant Time: Aug 30 11:57:48 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM082316.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 30 04:03:42 2016
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88 C	Benzo(a)pyrene	1.132	1.179	-4.2	73	0.00
89	Indeno(1,2,3-cd)pyrene	1.389	1.237	10.9	63	0.00
90	Dibenzo(a,h)anthracene	1.155	1.053	8.8	64	0.00
91	Benzo(g,h,i)perylene	1.179	0.997	15.4	60	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0