

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012317\
 Data File : BM008807.D
 Acq On : 23 Jan 2017 16:41
 Operator : UM/SJ
 Sample : SSTDICV020
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV41

Quant Time: Jan 23 17:18:03 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jan 23 16:57:33 2017
 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	104	0.00
2	1,4-Dioxane	0.490	0.582	-18.8	116	0.00
3 S	1,4-Dioxane-d8	0.405	0.452	-11.6	119	0.00
4	Benzaldehyde	1.394	1.732	-24.2#	106	0.00
5 S	Phenol-d5	1.694	1.764	-4.1	105	0.00
6	Phenol	1.847	1.955	-5.8	109	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.347	1.365	-1.3	100	0.00
8	Bis(2-Chloroethyl)ether	1.384	1.379	0.4	97	0.00
9 S	2-Chlorophenol-d4	1.104	1.156	-4.7	106	0.00
10	2-Chlorophenol	1.140	1.200	-5.3	104	0.00
11	2-Methylphenol	1.324	1.341	-1.3	104	0.00
12	2,2'-oxybis(1-Chloropropane	2.600	2.682	-3.2	100	0.00
13 S	4-Methylphenol-d8	1.326	1.367	-3.1	102	0.00
14	Acetophenone	2.431	2.593	-6.7	109	0.00
15 P	N-Nitroso-di-n-propylamine	1.584	1.618	-2.1	101	0.00
16	4-Methylphenol	1.427	1.416	0.8	100	0.00
17	Hexachloroethane	0.703	0.721	-2.6	106	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
19 S	Nitrobenzene-d5	0.127	0.140	-10.2	109	0.00
20	Nitrobenzene	0.561	0.605	-7.8	104	0.00
21	Isophorone	0.905	0.948	-4.8	101	0.00
22 S	2-Nitrophenol-d4	0.155	0.159	-2.6	101	0.00
23 C	2-Nitrophenol	0.160	0.172	-7.5	105	0.00
24	2,4-Dimethylphenol	0.458	0.486	-6.1	100	0.00
25	Bis(2-Chloroethoxy)methane	0.451	0.474	-5.1	100	0.00
26 S	2,4-Dichlorophenol-d3	0.324	0.331	-2.2	96	0.00
27 C	2,4-Dichlorophenol	0.318	0.326	-2.5	97	0.00
28	Naphthalene	0.923	0.964	-4.4	102	0.00
29 S	4-Chloroaniline-d4	0.353	0.378	-7.1	100	0.00
30	4-Chloroaniline	0.361	0.398	-10.2	105	0.00
31 C	Hexachlorobutadiene	0.308	0.308	0.0	95	0.00
32	Caprolactam	0.098	0.091	7.1	88	0.00
33 C	4-Chloro-3-methylphenol	0.413	0.432	-4.6	103	0.00
34	2-Methylnaphthalene	0.750	0.755	-0.7	96	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	92	0.00
36	1,2,4,5-Tetrachlorobenzene	0.614	0.662	-7.8	98	0.00
37	Hexachlorocyclopentadiene	0.475	0.490	-3.2	98	0.00
38 C	2,4,6-Trichlorophenol	0.371	0.392	-5.7	95	0.00
39	2,4,5-Trichlorophenol	0.398	0.419	-5.3	95	0.00
40	1,1'-Biphenyl	1.236	1.326	-7.3	96	0.00
41	2-Chloronaphthalene	0.972	1.059	-9.0	100	0.00
42	2-Nitroaniline	0.483	0.531	-9.9	97	0.00
43 S	Dimethylphthalate-d6	1.310	1.405	-7.3	95	0.00
44	Dimethylphthalate	1.310	1.409	-7.6	95	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012317\
 Data File : BM008807.D
 Acq On : 23 Jan 2017 16:41
 Operator : UM/SJ
 Sample : SSTDICV020
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV41

Quant Time: Jan 23 17:18:03 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jan 23 16:57:33 2017
 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	0.251	0.264	-5.2	92	0.00
46 S	Acenaphthylene-d8	1.550	1.658	-7.0	95	0.00
47	Acenaphthylene	1.494	1.615	-8.1	99	0.00
48	3-Nitroaniline	0.237	0.255	-7.6	88	0.00
49 C	Acenaphthene	1.051	1.128	-7.3	96	0.00
50	2,4-Dinitrophenol	0.180	0.164	8.9	87	0.00
51 S	4-Nitrophenol-d4	0.219	0.225	-2.7	91	0.00
52	4-Nitrophenol	0.452	0.464	-2.7	91	0.00
53	Dibenzofuran	1.616	1.713	-6.0	92	0.00
54	2,4-Dinitrotoluene	0.385	0.402	-4.4	91	0.00
55	2,3,4,6-Tetrachlorophenol	0.400	0.422	-5.5	91	0.00
56	Diethylphthalate	1.425	1.486	-4.3	92	0.00
57 S	Fluorene-d10	1.245	1.301	-4.5	91	0.00
58	Fluorene	1.350	1.408	-4.3	91	0.00
59	4-Chlorophenyl-phenylether	0.782	0.833	-6.5	94	0.00
60	4-Nitroaniline	0.265	0.281	-6.0	91	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.119	0.118	0.8	92	0.00
63	4,6-Dinitro-2-methylphenol	0.123	0.123	0.0	91	0.00
64	N-Nitrosodiphenylamine	0.506	0.541	-6.9	93	0.00
65	4-Bromophenyl-phenylether	0.217	0.231	-6.5	90	0.00
66	Hexachlorobenzene	0.236	0.250	-5.9	92	0.00
67	Atrazine	0.232	0.240	-3.4	88	0.00
68 C	Pentachlorophenol	0.155	0.153	1.3	89	0.00
69	Phenanthrene	0.985	1.040	-5.6	90	0.00
70 S	Anthracene-d10	0.875	0.929	-6.2	91	0.00
71	Anthracene	1.008	1.074	-6.5	90	0.00
72	Carbazole	0.884	0.909	-2.8	88	0.00
73	Di-n-butylphthalate	1.063	1.076	-1.2	85	0.00
74 C	Fluoranthene	1.280	1.290	-0.8	87	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	82	0.00
76 S	Pyrene-d10	0.877	0.945	-7.8	85	0.00
77	Pyrene	1.073	1.181	-10.1	88	0.00
78	Butylbenzylphthalate	0.400	0.420	-5.0	82	0.00
79	3,3'-Dichlorobenzidine	0.374	0.413	-10.4	82	0.00
80	Benzo(a)anthracene	1.075	1.134	-5.5	83	0.00
81	Bis(2-ethylhexyl)phthalate	0.555	0.573	-3.2	81	0.00
82	Chrysene	1.000	1.069	-6.9	84	0.00
83 I	Perylene-d12	1.000	1.000	0.0	83	0.00
84	Di-n-octyl phthalate	1.161	1.147	1.2	81	0.00
85	Benzo(b)fluoranthene	1.162	1.260	-8.4	82	0.00
86	Benzo(k)fluoranthene	1.104	1.155	-4.6	81	0.00
87 S	Benzo(a)pyrene-d12	0.847	0.896	-5.8	83	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012317\
Data File : BM008807.D
Acq On : 23 Jan 2017 16:41
Operator : UM/SJ
Sample : SSTDICV020
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SICV41

Quant Time: Jan 23 17:18:03 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jan 23 16:57:33 2017
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	1.066	1.138	-6.8	83	0.00
89	Indeno(1,2,3-cd)pyrene	1.070	1.136	-6.2	86	0.00
90	Dibenzo(a,h)anthracene	0.926	0.968	-4.5	85	0.00
91	Benzo(g,h,i)perylene	0.870	0.938	-7.8	86	0.00

(#) = Out of Range

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