

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020123\
 Data File : BP013688.D
 Acq On : 01 Feb 2023 10:42
 Operator : CG/JU
 Sample : SST01024
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SST010604

Quant Time: Feb 02 00:16:24 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 02 00:14:26 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.140	152	268542	20.000	ng/u1	0.00
20) Naphthalene-d8	10.969	136	1158459	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.775	164	870241	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.528	188	2081012	20.000	ng/u1	0.00
79) Chrysene-d12	21.604	240	2016635	20.000	ng/u1	0.00
88) Perylene-d12	24.174	264	2223534	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.481	96	26128	4.125	ng/uL	0.00
4) Pyridine-d5	3.917	84	181551	9.990	ng/u1	0.00
7) Phenol-d5	7.281	99	236077	10.003	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.458	67	145285	10.496	ng/u1	0.00
11) 2-Chlorophenol-d4	7.664	132	178260	10.303	ng/u1	0.00
15) 4-Methylphenol-d8	8.840	113	188911	10.040	ng/u1	0.00
21) Nitrobenzene-d5	9.310	128	66387	8.983	ng/u1	0.00
24) 2-Nitrophenol-d4	10.040	143	77049	8.946	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.575	165	199873	9.924	ng/u1	0.00
31) 4-Chloroaniline-d4	11.099	131	281520	10.584	ng/u1	0.00
46) Dimethylphthalate-d6	14.175	166	714012	10.409	ng/u1	0.00
49) Acenaphthylene-d8	14.469	160	776955	10.251	ng/u1	0.00
54) 4-Nitrophenol-d4	14.951	143	99493	8.605	ng/u1	0.00
60) Fluorene-d10	15.763	176	612763	10.363	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.875	200	81490	6.988	ng/u1	0.00
73) Anthracene-d10	17.628	188	1001779	10.327	ng/u1	0.00
81) Pyrene-d10	19.839	212	1163397	10.344	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.010	264	1151211	10.142	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.517	88	28099	4.257	ng/uL	97
5) Pyridine	3.934	79	187935	10.270	ng/u1	99
6) Benzaldehyde	7.269	77	95811	10.106	ng/u1	98
8) Phenol	7.311	94	245968	10.183	ng/u1	100
10) Bis(2-Chloroethyl)ether	7.546	93	201310	10.372	ng/u1	97
12) 2-Chlorophenol	7.699	128	179336	10.231	ng/u1	99
13) 2-Methylphenol	8.575	108	184457	10.002	ng/u1	98
14) 2,2'-oxybis(1-Chloropr...	8.663	45	247244	10.759	ng/u1	98
16) Acetophenone	8.969	105	318812	10.869	ng/u1	100
17) N-Nitroso-di-n-propyla...	8.946	70	156578	10.656	ng/u1	98
18) 4-Methylphenol	8.905	108	202930	10.132	ng/u1	100
19) Hexachloroethane	9.234	117	69287	10.099	ng/u1	99
22) Nitrobenzene	9.352	77	191775	9.608	ng/u1	100
23) Isophorone	9.881	82	471854	10.326	ng/u1	99
25) 2-Nitrophenol	10.069	139	89068	9.347	ng/u1	97
26) 2,4-Dimethylphenol	10.122	107	220493	10.241	ng/u1	97
27) Bis(2-Chloroethoxy)met...	10.363	93	288815	10.429	ng/u1	99
29) 2,4-Dichlorophenol	10.605	162	193056	9.905	ng/u1	98
30) Naphthalene	11.016	128	634932	10.370	ng/u1	99
32) 4-Chloroaniline	11.122	127	284125	10.613	ng/u1	99
33) Hexachlorobutadiene	11.305	225	138308	10.061	ng/u1	99
34) Caprolactam	11.893	113	66074	10.034	ng/u1	91
35) 4-Chloro-3-methylphenol	12.228	107	209938	10.162	ng/u1	99
36) 2-Methylnaphthalene	12.610	142	452231	10.376	ng/u1	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020123\
 Data File : BP013688.D
 Acq On : 01 Feb 2023 10:42
 Operator : CG/JU
 Sample : SSTD01024
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTD010604

Quant Time: Feb 02 00:16:24 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 02 00:14:26 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
37) 1-Methylnaphthalene	12.828	142	465146	10.567	ng/ul	100
39) 1,2,4,5-Tetrachloroben...	12.975	216	288023	10.019	ng/ul	99
40) Hexachlorocyclopentadiene	12.957	237	124169	7.683	ng/ul	99
41) 2,4,6-Trichlorophenol	13.210	196	187801	9.862	ng/ul	99
42) 2,4,5-Trichlorophenol	13.281	196	200010	9.638	ng/ul	99
43) 1,1'-Biphenyl	13.610	154	659227	10.005	ng/ul	99
44) 2-Chloronaphthalene	13.657	162	517864	9.934	ng/ul	99
45) 2-Nitroaniline	13.851	65	80484	8.111	ng/ul	96
47) Dimethylphthalate	14.222	163	709198	10.494	ng/ul	99
48) 2,6-Dinitrotoluene	14.340	165	76609	7.954	ng/ul#	94
50) Acenaphthylene	14.498	152	829781	10.383	ng/ul	99
51) 3-Nitroaniline	14.675	138	82138	8.069	ng/ul	98
52) Acenaphthene	14.840	153	576904	10.504	ng/ul	99
53) 2,4-Dinitrophenol	14.875	184	45387	6.067	ng/ul	98
55) 4-Nitrophenol	14.963	109	78563	8.913	ng/ul	95
56) Dibenzofuran	15.169	168	848665	10.638	ng/ul	99
57) 2,4-Dinitrotoluene	15.128	165	129640	8.570	ng/ul#	99
58) 2,3,4,6-Tetrachlorophenol	15.393	232	195385	10.017	ng/ul	99
59) Diethylphthalate	15.581	149	674717	10.400	ng/ul	99
61) Fluorene	15.822	166	690133	10.630	ng/ul	99
62) 4-Chlorophenyl-phenyle...	15.810	204	372761	10.539	ng/ul	99
63) 4-Nitroaniline	15.834	138	71286	7.479	ng/ul	99
66) 4,6-Dinitro-2-methylph...	15.887	198	85533	7.370	ng/ul#	95
67) N-Nitrosodiphenylamine	16.022	169	592541	10.549	ng/ul	99
68) 4-Bromophenyl-phenylether	16.710	248	240700	10.166	ng/ul	98
69) Hexachlorobenzene	16.828	284	285668	10.284	ng/ul	99
70) Atrazine	16.969	200	222588	9.846	ng/ul	99
71) Pentachlorophenol	17.169	266	155960	8.797	ng/ul	98
72) Phenanthrene	17.569	178	1145091	10.417	ng/ul	99
74) Anthracene	17.663	178	1162997	10.446	ng/ul	100
75) 1,2,3,4-Tetrachloroben...	13.581	216	283131	9.550	ng/uL	99
76) Pentachlorobenzene	15.093	250	312854	10.293	ng/uL	97
77) Carbazole	17.922	167	1032393	10.689	ng/ul	99
78) Di-n-butylphthalate	18.457	149	1144165	10.026	ng/ul	99
80) Fluoranthene	19.516	202	1413218	10.780	ng/ul	97
82) Pyrene	19.869	202	1411903	10.534	ng/ul	97
83) Butylbenzylphthalate	20.728	149	380087	9.870	ng/ul	99
84) 3,3'-Dichlorobenzidine	21.510	252	416635	9.096	ng/ul	100
85) Benzo(a)anthracene	21.586	228	1438864	10.378	ng/ul	99
86) Bis(2-ethylhexyl)phtha...	21.486	149	659886	9.995	ng/ul	99
87) Chrysene	21.645	228	1360313	10.422	ng/ul	98
89) Di-n-octyl phthalate	22.469	149	1139317	9.727	ng/ul	100
90) Benzo(b)fluoranthene	23.386	252	1486312	10.240	ng/ul	99
91) Benzo(k)fluoranthene	23.433	252	1423048	10.258	ng/ul	99
93) Benzo(a)pyrene	24.063	252	1280141	10.216	ng/ul	100
94) Indeno(1,2,3-cd)pyrene	26.892	276	1570093	10.113	ng/ul	98
95) Dibenzo(a,h)anthracene	26.910	278	1295448	10.065	ng/ul	99
96) Benzo(g,h,i)perylene	27.727	276	1230299	10.017	ng/ul	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

