

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100122\
 Data File : BP011852.D
 Acq On : 01 Oct 2022 09:36
 Operator : CG/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTD020760

Quant Time: Oct 03 00:57:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 01 00:53:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	241180	20.000	ng/u1	0.00
20) Naphthalene-d8	10.704	136	1129773	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.539	164	752243	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.298	188	1646567	20.000	ng/u1	0.00
79) Chrysene-d12	21.386	240	1708482	20.000	ng/u1	0.00
88) Perylene-d12	23.821	264	1864614	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	48623	8.891	ng/uL	0.00
4) Pyridine-d5	3.728	84	345491	21.569	ng/u1	0.00
7) Phenol-d5	7.052	99	444204	21.334	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	256603	22.234	ng/u1	0.00
11) 2-Chlorophenol-d4	7.422	132	362803	22.073	ng/u1	0.00
15) 4-Methylphenol-d8	8.604	113	375660	21.826	ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	186000	21.614	ng/u1	0.00
24) 2-Nitrophenol-d4	9.781	143	202567	22.002	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	381468	22.401	ng/u1	0.00
31) 4-Chloroaniline-d4	10.840	131	582079	22.421	ng/u1	0.00
46) Dimethylphthalate-d6	13.951	166	1210365	22.500	ng/u1	0.00
49) Acenaphthylene-d8	14.234	160	1386072	22.602	ng/u1	0.00
54) 4-Nitrophenol-d4	14.734	143	244603	21.958	ng/u1	0.00
60) Fluorene-d10	15.534	176	1042435	22.314	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	210563	21.128	ng/u1	0.00
73) Anthracene-d10	17.398	188	1662009	22.159	ng/u1	0.00
81) Pyrene-d10	19.633	212	1909108	22.064	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.663	264	2077584	22.258	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.340	88	50868	8.612	ng/uL	93
5) Pyridine	3.746	79	341612	21.973	ng/u1	99
6) Benzaldehyde	7.034	77	234608	25.722	ng/u1	99
8) Phenol	7.081	94	464132	21.492	ng/u1	99
10) Bis(2-Chloroethyl)ether	7.316	93	380970	22.170	ng/u1	99
12) 2-Chlorophenol	7.452	128	389768	22.182	ng/u1	99
13) 2-Methylphenol	8.334	108	367667	21.609	ng/u1	98
14) 2,2'-oxybis(1-Chloropr...	8.428	45	378837	22.363	ng/u1	98
16) Acetophenone	8.722	105	595684	22.700	ng/u1	99
17) N-Nitroso-di-n-propyla...	8.704	70	283038	22.916	ng/u1	99
18) 4-Methylphenol	8.663	108	410689	21.908	ng/u1	98
19) Hexachloroethane	8.975	117	142871	21.800	ng/u1	98
22) Nitrobenzene	9.104	77	409564	22.130	ng/u1	100
23) Isophorone	9.634	82	843084	22.456	ng/u1	100
25) 2-Nitrophenol	9.816	139	232402	22.101	ng/u1	98
26) 2,4-Dimethylphenol	9.875	107	440700	22.310	ng/u1	98
27) Bis(2-Chloroethoxy)met...	10.116	93	542101	22.380	ng/u1	99
29) 2,4-Dichlorophenol	10.346	162	395423	22.410	ng/u1	98
30) Naphthalene	10.751	128	1330161	21.920	ng/u1	100
32) 4-Chloroaniline	10.863	127	603511	22.722	ng/u1	99
33) Hexachlorobutadiene	11.040	225	227920	21.822	ng/u1	98
34) Caprolactam	11.651	113	131971	21.133	ng/u1	94
35) 4-Chloro-3-methylphenol	11.987	107	424589	22.694	ng/u1	99
36) 2-Methylnaphthalene	12.363	142	944876	22.297	ng/u1	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
37) 1-Methylnaphthalene	12.581	142	949472	22.329	ng/u1	98
39) 1,2,4,5-Tetrachloroben...	12.728	216	488790	22.406	ng/u1	98
40) Hexachlorocyclopentadiene	12.710	237	261112	20.952	ng/u1	99
41) 2,4,6-Trichlorophenol	12.969	196	324830	22.244	ng/u1	99
42) 2,4,5-Trichlorophenol	13.040	196	358471	22.535	ng/u1	98
43) 1,1'-Biphenyl	13.375	154	1250655	22.478	ng/u1	100
44) 2-Chloronaphthalene	13.416	162	984218	22.374	ng/u1	99
45) 2-Nitroaniline	13.622	65	241907	22.528	ng/u1	98
47) Dimethylphthalate	13.998	163	1275491	22.625	ng/u1	99
48) 2,6-Dinitrotoluene	14.116	165	249781	22.734	ng/u1	97
50) Acenaphthylene	14.263	152	1544265	22.674	ng/u1	99
51) 3-Nitroaniline	14.445	138	275735	21.652	ng/u1	99
52) Acenaphthene	14.604	153	1078733	22.449	ng/u1	99
53) 2,4-Dinitrophenol	14.651	184	147096	19.747	ng/u1	97
55) 4-Nitrophenol	14.751	109	167761	22.309	ng/u1	98
56) Dibenzofuran	14.939	168	1488646	22.618	ng/u1	97
57) 2,4-Dinitrotoluene	14.904	165	370590	23.076	ng/u1	100
58) 2,3,4,6-Tetrachlorophenol	15.163	232	311876	22.569	ng/u1	100
59) Diethylphthalate	15.363	149	1276030	22.593	ng/u1	100
61) Fluorene	15.592	166	1220569	22.643	ng/u1	98
62) 4-Chlorophenyl-phenyle...	15.586	204	603594	22.618	ng/u1	99
63) 4-Nitroaniline	15.616	138	271775	22.365	ng/u1	99
66) 4,6-Dinitro-2-methylph...	15.669	198	227018	21.502	ng/u1	98
67) N-Nitrosodiphenylamine	15.798	169	1070037	22.030	ng/u1	99
68) 4-Bromophenyl-phenylether	16.486	248	364707	21.776	ng/u1	98
69) Hexachlorobenzene	16.598	284	418014	21.933	ng/u1	99
70) Atrazine	16.757	200	385643	22.004	ng/u1	99
71) Pentachlorophenol	16.945	266	261814	20.886	ng/u1	99
72) Phenanthrene	17.339	178	2003366	22.139	ng/u1	100
74) Anthracene	17.433	178	2030961	22.307	ng/u1	99
75) 1,2,3,4-Tetrachloroben...	13.340	216	489262	21.528	ng/uL	99
76) Pentachlorobenzene	14.857	250	531110	21.870	ng/uL	99
77) Carbazole	17.704	167	1866401	22.689	ng/u1	99
78) Di-n-butylphthalate	18.251	149	2223253	22.070	ng/u1	100
80) Fluoranthene	19.304	202	2335490	22.056	ng/u1	100
82) Pyrene	19.657	202	2450666	22.228	ng/u1	99
83) Butylbenzylphthalate	20.533	149	1041856	21.865	ng/u1	100
84) 3,3'-Dichlorobenzidine	21.304	252	912139	23.191	ng/u1	100
85) Benzo(a)anthracene	21.369	228	2504489	22.416	ng/u1	100
86) Bis(2-ethylhexyl)phtha...	21.286	149	1565848	21.816	ng/u1	99
87) Chrysene	21.421	228	2334864	22.268	ng/u1	99
89) Di-n-octyl phthalate	22.227	149	2708036	22.764	ng/u1	100
90) Benzo(b)fluoranthene	23.074	252	2620518	22.019	ng/u1	99
91) Benzo(k)fluoranthene	23.127	252	2637597	23.014	ng/u1	99
93) Benzo(a)pyrene	23.715	252	2377649	22.387	ng/u1	99
94) Indeno(1,2,3-cd)pyrene	26.362	276	2949516	21.818	ng/u1	100
95) Dibenzo(a,h)anthracene	26.380	278	2653855	22.099	ng/u1	99
96) Benzo(g,h,i)perylene	27.145	276	2603432	21.722	ng/u1	98

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

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