

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010852.D
 Acq On : 26 Jun 2022 05:08
 Operator : CG/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTD020683

Quant Time: Jun 26 22:56:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 24 22:42:26 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 06/27/2022
 Supervised By :mohammad ahmed 07/01/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	904813	20.000	ng/u1	0.00
20) Naphthalene-d8	10.510	136	4107988	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.375	164	2318934	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.134	188	4601333	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.251	240	4537929	20.000	ng/u1	# 0.00
88) Perylene-d12	23.616	264	5020284	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	171049	7.151	ng/uL	0.00
4) Pyridine-d5	3.611	84	1248558	19.610	ng/u1	0.00
7) Phenol-d5	6.893	99	1517456	21.186	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.058	67	949018	20.260	ng/u1	0.00
11) 2-Chlorophenol-d4	7.252	132	1245581	20.961	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	1253375	22.136	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	599774	22.935	ng/u1	0.00
24) 2-Nitrophenol-d4	9.599	143	597794	27.630	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.134	165	1094363	20.471	ng/u1	0.00
31) 4-Chloroaniline-d4	10.657	131	1708756	19.582	ng/u1	0.00
46) Dimethylphthalate-d6	13.793	166	2987203	18.805	ng/u1	0.00
49) Acenaphthylene-d8	14.063	160	3746291	18.479	ng/u1	0.00
54) 4-Nitrophenol-d4	14.592	143	608215	24.514	ng/u1	0.01
60) Fluorene-d10	15.369	176	2554685	18.741	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.492	200	382075	25.063	ng/u1	0.00
73) Anthracene-d10	17.234	188	3811367	18.617	ng/u1	-0.01
81) Pyrene-d10	19.486	212	4375854	17.700	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.463	264	4692342	18.915	ng/u1	-0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.240	88	177991	6.978	ng/uL#	84
5) Pyridine	3.628	79	1265829	19.626	ng/u1#	73
6) Benzaldehyde	6.864	77	568106	15.032	ng/u1	94
8) Phenol	6.922	94	1608207	20.993	ng/u1#	85
10) Bis(2-Chloroethyl)ether	7.152	93	1269286	20.630	ng/u1	88
12) 2-Chlorophenol	7.281	128	1271890	20.735	ng/u1	92
13) 2-Methylphenol	8.169	108	1206307	21.152	ng/u1	99
14) 2,2'-oxybis(1-Chloropr...	8.246	45	1665327	20.174	ng/u1#	97
16) Acetophenone	8.546	105	1850120	21.277	ng/u1#	87
17) N-Nitroso-di-n-propyla...	8.540	70	911656	20.734	ng/u1#	87
18) 4-Methylphenol	8.499	108	1294158	21.790	ng/u1	97
19) Hexachloroethane	8.787	117	487148	19.610	ng/u1#	82
22) Nitrobenzene	8.922	77	1337749	20.498	ng/u1#	88
23) Isophorone	9.452	82	2534272	18.517	ng/u1	97
25) 2-Nitrophenol	9.628	139	678302	25.889	ng/u1#	78
26) 2,4-Dimethylphenol	9.699	107	1354184	18.918	ng/u1	93
27) Bis(2-Chloroethoxy)met...	9.940	93	1635706	18.946	ng/u1	97
29) 2,4-Dichlorophenol	10.163	162	1111792	20.015	ng/u1	99
30) Naphthalene	10.563	128	3922839	18.322	ng/u1	99
32) 4-Chloroaniline	10.681	127	1698821	19.649	ng/u1	100
33) Hexachlorobutadiene	10.846	225	631558	17.796	ng/u1	95
34) Caprolactam	11.475	113	381864m	21.883	ng/u1	
35) 4-Chloro-3-methylphenol	11.816	107	1231392	20.453	ng/u1	93
36) 2-Methylnaphthalene	12.181	142	2605838	18.795	ng/u1	99

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37) 1-Methylnaphthalene	12.404	142	2609882	18.948	ng/ul#	99
39) 1,2,4,5-Tetrachloroben...	12.551	216	1203651	18.063	ng/ul	99
40) Hexachlorocyclopentadiene	12.528	237	774130	18.333	ng/ul	96
41) 2,4,6-Trichlorophenol	12.798	196	805172	20.958	ng/ul	98
42) 2,4,5-Trichlorophenol	12.869	196	883046	21.419	ng/ul	98
43) 1,1'-Biphenyl	13.204	154	3328802	18.235	ng/ul#	97
44) 2-Chloronaphthalene	13.240	162	2511454	18.263	ng/ul	98
45) 2-Nitroaniline	13.457	65	739981	25.362	ng/ul#	76
47) Dimethylphthalate	13.840	163	2966132	18.645	ng/ul	98
48) 2,6-Dinitrotoluene	13.957	165	611610	25.660	ng/ul	90
50) Acenaphthylene	14.093	152	4089748	18.420	ng/ul	99
51) 3-Nitroaniline	14.287	138	609453	21.328	ng/ul	88
52) Acenaphthene	14.434	153	2621867	18.301	ng/ul	99
53) 2,4-Dinitrophenol	14.487	184	290475	31.744	ng/ul#	79
55) 4-Nitrophenol	14.604	109	453764	23.000	ng/ul#	76
56) Dibenzofuran	14.775	168	3644269	18.571	ng/ul	95
57) 2,4-Dinitrotoluene	14.745	165	862783	27.090	ng/ul#	78
58) 2,3,4,6-Tetrachlorophenol	15.004	232	660229	22.363	ng/ul	94
59) Diethylphthalate	15.216	149	2999677	18.967	ng/ul	99
61) Fluorene	15.428	166	2894698	18.499	ng/ul	99
62) 4-Chlorophenyl-phenyle...	15.428	204	1355053	18.499	ng/ul	99
63) 4-Nitroaniline	15.457	138	574731	21.753	ng/ul#	76
66) 4,6-Dinitro-2-methylph...	15.510	198	398881	24.214	ng/ul#	93
67) N-Nitrosodiphenylamine	15.645	169	2504531	17.971	ng/ul	99
68) 4-Bromophenyl-phenylether	16.328	248	830232	18.004	ng/ul	97
69) Hexachlorobenzene	16.428	284	952681	17.836	ng/ul	96
70) Atrazine	16.610	200	854486	18.493	ng/ul	97
71) Pentachlorophenol	16.781	266	676386	24.593	ng/ul	98
72) Phenanthrene	17.181	178	4504733	18.384	ng/ul	99
74) Anthracene	17.269	178	4530109	18.382	ng/ul	99
75) 1,2,3,4-Tetrachloroben...	13.163	216	1253785	16.512	ng/uL	98
76) Pentachlorobenzene	14.687	250	1133977	17.413	ng/uL	100
77) Carbazole	17.545	167	4302950	19.397	ng/ul	99
78) Di-n-butylphthalate	18.116	149	5068211	19.172	ng/ul	99
80) Fluoranthene	19.157	202	5322106	17.503	ng/ul#	87
82) Pyrene	19.516	202	5357479	17.447	ng/ul#	85
83) Butylbenzylphthalate	20.410	149	2414052	22.335	ng/ul	94
84) 3,3'-Dichlorobenzidine	21.175	252	1683610	17.967	ng/ul	98
85) Benzo(a)anthracene	21.233	228	5304427	18.690	ng/ul	97
86) Bis(2-ethylhexyl)phtha...	21.175	149	3467231	20.257	ng/ul#	97
87) Chrysene	21.286	228	5049506	18.567	ng/ul	99
89) Di-n-octyl phthalate	22.098	149	6002173	20.142	ng/ul	100
90) Benzo(b)fluoranthene	22.892	252	5818321	18.637	ng/ul#	96
91) Benzo(k)fluoranthene	22.945	252	5481702	18.472	ng/ul#	95
93) Benzo(a)pyrene	23.516	252	5657411	18.628	ng/ul#	96
94) Indeno(1,2,3-cd)pyrene	26.074	276	6803722	18.566	ng/ul#	90
95) Dibenzo(a,h)anthracene	26.098	278	5803821	18.463	ng/ul#	92
96) Benzo(g,h,i)perylene	26.833	276	5710245	18.318	ng/ul#	88

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

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