

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070823\
 Data File : BP016132.D
 Acq On : 09 Jul 2023 14:32
 Operator : MA/JU
 Sample : 03356-10MS
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YBW38MS

Quant Time: Jul 09 22:41:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 08 10:11:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.011	152	290233	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.828	136	1185384	20.000	ng/u1	#-0.02
38) Acenaphthene-d10	14.663	164	676662	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.428	188	1288448	20.000	ng/u1	#-0.02
79) Chrysene-d12	21.527	240	923238	20.000	ng/u1	#-0.01
88) Perylene-d12	24.074	264	1083179	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	44309	5.493	ng/uL	0.00
4) Pyridine-d5	3.793	84	524755	25.806	ng/u1	-0.01
7) Phenol-d5	7.193	99	806223	32.466	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.340	67	525717	34.219	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.540	132	600627	32.041	ng/u1	-0.02
15) 4-Methylphenol-d8	8.752	113	618009	31.503	ng/u1	-0.02
21) Nitrobenzene-d5	9.199	128	296239	32.701	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.928	143	333127	31.507	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.481	165	589224	31.273	ng/u1	-0.02
31) 4-Chloroaniline-d4	11.004	131	734193	30.334	ng/u1	-0.02
46) Dimethylphthalate-d6	14.075	166	1610172	30.787	ng/u1	-0.01
49) Acenaphthylene-d8	14.357	160	1896100	32.250	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.951	143	71686	10.387	ng/u1	-0.01
60) Fluorene-d10	15.657	176	1330379	30.893	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.839	200	191651	24.187	ng/u1	-0.01
73) Anthracene-d10	17.534	188	1862264	31.642	ng/u1	-0.01
81) Pyrene-d10	19.763	212	1899457	31.508	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.898	264	1765821	32.317	ng/u1	-0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.381	88	107725	12.409	ng/uL	97
5) Pyridine	3.811	79	647366	30.431	ng/u1	97
6) Benzaldehyde	7.163	77	470513	46.864	ng/u1	97
8) Phenol	7.222	94	859543	33.355	ng/u1	93
10) Bis(2-Chloroethyl)ether	7.434	93	711110	34.683	ng/u1	98
12) 2-Chlorophenol	7.575	128	645843	32.429	ng/u1	98
13) 2-Methylphenol	8.475	108	633749	32.573	ng/u1	98
14) 2,2'-oxybis(1-Chloropr...	8.546	45	1072368	38.293	ng/u1	98
16) Acetophenone	8.858	105	1031828	31.996	ng/u1	97
17) N-Nitroso-di-n-propyla...	8.834	70	571257	31.895	ng/u1	98
18) 4-Methylphenol	8.822	108	691661	33.086	ng/u1	94
19) Hexachloroethane	9.075	117	287857	31.295	ng/u1	89
22) Nitrobenzene	9.246	77	854819	33.601	ng/u1	98
23) Isophorone	9.769	82	1575921	32.371	ng/u1	99
25) 2-Nitrophenol	9.957	139	371140	33.075	ng/u1	93
26) 2,4-Dimethylphenol	10.016	107	759852	30.804	ng/u1	97
27) Bis(2-Chloroethoxy)met...	10.240	93	955334	34.762	ng/u1	98
29) 2,4-Dichlorophenol	10.510	162	597767	31.912	ng/u1	99
30) Naphthalene	10.881	128	2069086	32.109	ng/u1	99
32) 4-Chloroaniline	11.028	127	671921	26.720	ng/u1	98
33) Hexachlorobutadiene	11.140	225	388471	27.462	ng/u1	99
34) Caprolactam	11.846	113	179141	31.265	ng/u1	94
35) 4-Chloro-3-methylphenol	12.169	107	665758	30.364	ng/u1	99
36) 2-Methylnaphthalene	12.493	142	1344782	31.701	ng/u1	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
37) 1-Methylnaphthalene	12.710	142	1334301	30.833	ng/ul	100
39) 1,2,4,5-Tetrachloroben...	12.863	216	704990	31.616	ng/ul	99
40) Hexachlorocyclopentadiene	12.810	237	207652	31.128	ng/ul	97
41) 2,4,6-Trichlorophenol	13.122	196	440023	31.225	ng/ul	99
42) 2,4,5-Trichlorophenol	13.216	196	475690	31.825	ng/ul	97
43) 1,1'-Biphenyl	13.493	154	1774998	33.118	ng/ul#	97
44) 2-Chloronaphthalene	13.546	162	1385616	32.818	ng/ul	98
45) 2-Nitroaniline	13.787	65	450405	33.903	ng/ul	97
47) Dimethylphthalate	14.122	163	1660767	30.683	ng/ul	99
48) 2,6-Dinitrotoluene	14.275	165	344227	31.352	ng/ul	100
50) Acenaphthylene	14.387	152	2107540	32.533	ng/ul	100
51) 3-Nitroaniline	14.628	138	318467	36.989	ng/ul	95
52) Acenaphthene	14.728	153	1458522	32.468	ng/ul	99
53) 2,4-Dinitrophenol	14.875	184	145465	24.972	ng/ul	88
55) 4-Nitrophenol	14.981	109	224941	31.450	ng/ul#	51
56) Dibenzofuran	15.069	168	1946893	31.220	ng/ul	97
57) 2,4-Dinitrotoluene	15.087	165	462870	30.937	ng/ul	86
58) 2,3,4,6-Tetrachlorophenol	15.316	232	370339	28.845	ng/ul	98
59) Diethylphthalate	15.475	149	1646602	30.062	ng/ul	99
61) Fluorene	15.716	166	1563208	30.905	ng/ul	98
62) 4-Chlorophenyl-phenyle...	15.698	204	760237	29.373	ng/ul	98
63) 4-Nitroaniline	15.804	138	262556	44.556	ng/ul	86
66) 4,6-Dinitro-2-methylph...	15.857	198	239693	29.897	ng/ul#	93
67) N-Nitrosodiphenylamine	15.928	169	1271760	32.971	ng/ul	99
68) 4-Bromophenyl-phenylether	16.604	248	461664	31.378	ng/ul	90
69) Hexachlorobenzene	16.728	284	515488	29.693	ng/ul	97
70) Atrazine	16.892	200	322891	21.869	ng/ul	96
71) Pentachlorophenol	17.098	266	213779	25.657	ng/ul	97
72) Phenanthrene	17.469	178	2265145	32.361	ng/ul	100
74) Anthracene	17.569	178	2243815	31.902	ng/ul	99
75) 1,2,3,4-Tetrachloroben...	13.469	216	717131	34.205	ng/ul	100
76) Pentachlorobenzene	14.987	250	1504389	71.097	ng/ul	99
77) Carbazole	17.863	167	1892842	31.930	ng/ul	99
78) Di-n-butylphthalate	18.357	149	2608490	33.191	ng/ul	100
80) Fluoranthene	19.439	202	2331402	31.117	ng/ul#	94
82) Pyrene	19.798	202	2384438	31.199	ng/ul#	91
83) Butylbenzylphthalate	20.639	149	1004712	32.223	ng/ul	94
84) 3,3'-Dichlorobenzidine	21.451	252	611796	29.314	ng/ul	97
85) Benzo(a)anthracene	21.510	228	2034259	31.323	ng/ul	99
86) Bis(2-ethylhexyl)phtha...	21.380	149	1470140	34.308	ng/ul	100
87) Chrysene	21.569	228	2053396	33.325	ng/ul	99
89) Di-n-octyl phthalate	22.333	149	2450621	30.958	ng/ul	100
90) Benzo(b)fluoranthene	23.274	252	2118428	30.438	ng/ul	98
91) Benzo(k)fluoranthene	23.333	252	2098938	29.848	ng/ul#	98
93) Benzo(a)pyrene	23.957	252	1928865	31.717	ng/ul#	98
94) Indeno(1,2,3-cd)pyrene	26.745	276	2631929	31.880	ng/ul#	93
95) Dibenzo(a,h)anthracene	26.745	278	2154158	31.767	ng/ul#	95
96) Benzo(g,h,i)perylene	27.598	276	2166982	31.906	ng/ul#	94

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

