

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121619\
 Data File : BM024114.D
 Acq On : 16 Dec 2019 18:33
 Operator : JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

mohammad
 12/17/2019 3:20:23 PM

Quant Time: Dec 16 19:09:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 16 17:25:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	390642	20.00	ng	0.00
21) Naphthalene-d8	10.23	136	1627889	20.00	ng	0.00
39) Acenaphthene-d10	14.12	164	1072076	20.00	ng	0.00
64) Phenanthrene-d10	16.87	188	2328791	20.00	ng	0.00
76) Chrysene-d12	21.09	240	2316640	20.00	ng	0.00
87) Perylene-d12	23.22	264	2527428	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.08	112	3456200	140.32	ng	0.00
7) Phenol-d6	6.66	99	5338872	168.76	ng	0.01
23) Nitrobenzene-d5	8.62	82	5571949	190.00	ng	0.00
42) 2,4,6-Tribromophenol	15.63	330	2608409	194.90	ng	0.00
45) 2-Fluorobiphenyl	12.73	172	12470016	160.17	ng	0.00
79) Terphenyl-d14	19.52	244	14995903	125.34	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.03	88	709660	71.602	ng	# 100
3) Pyridine	3.43	79	1999540	76.540	ng	98
4) n-Nitrosodimethylamine	3.35	42	697830	73.376	ng	93
6) Aniline	6.80	93	3355748	88.464	ng	100
8) 2-Chlorophenol	7.03	128	2034879	81.084	ng	99
9) Benzaldehyde	6.62	77	1019209	56.580	ng	98
10) Phenol	6.69	94	2777252	93.247	ng	99
11) bis(2-Chloroethyl)ether	6.90	93	2119886	90.593	ng	98
12) 1,3-Dichlorobenzene	7.35	146	2269438	75.662	ng	99
13) 1,4-Dichlorobenzene	7.49	146	2311087	76.109	ng	98
14) 1,2-Dichlorobenzene	7.80	146	2272792	78.509	ng	99
15) Benzyl Alcohol	7.71	79	2109168	106.547	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.00	45	2525112	74.286	ng	98
17) 2-Methylphenol	7.92	107	1870128	94.863	ng	98
18) Hexachloroethane	8.52	117	873621	78.442	ng	98
19) n-Nitroso-di-n-propylamine	8.28	70	1927106	109.085	ng	97
20) 3+4-Methylphenols	8.25	107	2582986	98.370	ng	99
22) Acetophenone	8.29	105	3558144	88.549	ng	99
24) Nitrobenzene	8.66	77	2709197	96.554	ng	98
25) Isophorone	9.19	82	5088289	105.524	ng	100
26) 2-Nitrophenol	9.36	139	1258934	103.932	ng	97
27) 2,4-Dimethylphenol	9.43	122	1767515	87.162	ng	100
28) bis(2-Chloroethoxy)methane	9.66	93	3160515	94.719	ng	98
29) 2,4-Dichlorophenol	9.89	162	2111182	92.084	ng	100
30) 1,2,4-Trichlorobenzene	10.10	180	2351549	83.366	ng	98
31) Naphthalene	10.27	128	6982935	86.815	ng	99
32) Benzoic acid	9.67	122	1711866m	131.162	ng	
33) 4-Chloroaniline	10.40	127	3177129	90.974	ng	98
34) Hexachlorobutadiene	10.56	225	1353923	80.520	ng	98
35) Caprolactam	11.25	113	805543m	118.516	ng	
36) 4-Chloro-3-methylphenol	11.55	107	2494625	111.525	ng	98
37) 2-Methylnaphthalene	11.90	142	5200738	92.889	ng	99
38) 1-Methylnaphthalene	12.13	142	4945321	93.052	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.29	216	2631114	72.430	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.26	237	1091093	55.050	nq	99
43) 2,4,6-Trichlorophenol	12.54	196	1783222	85.052	nq	100
44) 2,4,5-Trichlorophenol	12.62	196	1832621	83.879	nq	97
46) 1,1'-Biphenyl	12.95	154	6918068	77.969	nq	99
47) 2-Chloronaphthalene	12.98	162	5482588	82.902	nq	100
48) 2-Nitroaniline	13.20	65	1891030	109.811	nq	97
49) Acenaphthylene	13.84	152	8536055	87.162	nq	100
50) Dimethylphthalate	13.60	163	7001354	90.368	nq	100
51) 2,6-Dinitrotoluene	13.72	165	1566534	96.664	nq	91
52) Acenaphthene	14.19	154	5215797	81.407	nq	99
53) 3-Nitroaniline	14.05	138	1781550	95.491	nq	95
54) 2,4-Dinitrophenol	14.27	184	965228	126.144	nq	97
55) Dibenzofuran	14.53	168	8196051	87.624	nq	99
56) 4-Nitrophenol	14.38	139	1335971	93.164	nq	98
57) 2,4-Dinitrotoluene	14.52	165	2237513	105.390	nq	97
58) Fluorene	15.18	166	6855217	90.189	nq	100
59) 2,3,4,6-Tetrachlorophenol	14.76	232	1654439	86.827	nq	99
60) Diethylphthalate	14.97	149	7113726	94.430	nq	99
61) 4-Chlorophenyl-phenylether	15.18	204	3435255	87.960	nq	99
62) 4-Nitroaniline	15.23	138	1767818	90.380	nq	99
63) Azobenzene	15.47	77	6852790	100.731	nq	99
65) 4,6-Dinitro-2-methylphenol	15.29	198	1213434	111.260	nq	99
66) n-Nitrosodiphenylamine	15.40	169	5926095	83.297	nq	100
67) 4-Bromophenyl-phenylether	16.07	248	2280679	88.051	nq	98
68) Hexachlorobenzene	16.19	284	2735569	90.563	nq	99
69) Atrazine	16.36	200	901791	38.992	nq	98
70) Pentachlorophenol	16.54	266	1321204	82.016	nq	100
71) Phenanthrene	16.92	178	10749837	81.808	nq	99
72) Anthracene	17.01	178	10511957	82.543	nq	99
73) Carbazole	17.29	167	9935729	83.841	nq	99
74) Di-n-butylphthalate	17.86	149	11444552	85.169	nq	96
75) Fluoranthene	18.94	202	12149528	81.937	nq	95
77) Benzidine	19.14	184	2779002	43.727	nq	99
78) Pyrene	19.31	202	12261264	80.762	nq	95
80) Butylbenzylphthalate	20.24	149	5970221	103.471	nq	95
81) Benzo(a)anthracene	21.07	228	11882292	79.139	nq	91
82) 3,3'-Dichlorobenzidine	21.02	252	4391114	82.468	nq	98
83) Chrysene	21.13	228	11238649m	77.209	nq	
84) Bis(2-ethylhexyl)phthalate	21.03	149	8352304	97.696	nq	# 96
85) Di-n-octyl phthalate	21.87	149	12157047	89.222	nq	98
86) Indeno(1,2,3-cd)pyrene	25.35	276	13173732	75.826	nq	# 100
88) Benzo(b)fluoranthene	22.60	252	13427907	87.748	nq	99
89) Benzo(k)fluoranthene	22.64	252	12391412	81.609	nq	97
90) Benzo(a)pyrene	23.13	252	11986667	85.006	nq	99
91) Dibenzo(a,h)anthracene	25.35	278	11004062	74.606	nq	99
92) Benzo(g,h,i)perylene	25.99	276	10040882	71.673	nq	100

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

