

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090220\
 Data File : BM027306.D
 Acq On : 02 Sep 2020 16:37
 Operator : JU/CG
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICBM090220

Quant Time: Sep 02 17:48:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM090220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 16:21:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	150227	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	585419	20.00	ng	0.00
39) Acenaphthene-d10	14.22	164	421993	20.00	ng	0.00
64) Phenanthrene-d10	16.97	188	983505	20.00	ng	0.00
76) Chrysene-d12	21.17	240	1093096	20.00	ng	0.00
86) Perylene-d12	23.34	264	1098820	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.20	112	704914	78.12	ng	0.00
7) Phenol-d6	6.76	99	986488	81.01	ng	0.00
23) Nitrobenzene-d5	8.72	82	1055642	77.04	ng	0.00
42) 2,4,6-Tribromophenol	15.72	330	581455	78.58	ng	0.00
45) 2-Fluorobiphenyl	12.84	172	2459449	77.51	ng	0.00
79) Terphenyl-d14	19.62	244	4780719	76.12	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.17	88	144871	37.655	ng	99
3) Pyridine	3.55	79	438415	39.227	ng	99
4) n-Nitrosodimethylamine	3.46	42	165830	37.596	ng	96
6) Aniline	6.92	93	586168	39.603	ng	98
8) 2-Chlorophenol	7.15	128	361911	39.399	ng	98
9) Benzaldehyde	6.73	77	374983	39.143	ng	97
10) Phenol	6.79	94	468773	40.063	ng	97
11) bis(2-Chloroethyl)ether	7.01	93	392130	39.824	ng	99
12) 1,3-Dichlorobenzene	7.47	146	443579	38.795	ng	98
13) 1,4-Dichlorobenzene	7.61	146	446707	38.856	ng	99
14) 1,2-Dichlorobenzene	7.93	146	427029	38.655	ng	99
15) Benzyl Alcohol	7.82	79	419085	40.048	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.12	45	300472	40.295	ng	97
17) 2-Methylphenol	8.02	107	320950	40.394	ng	98
18) Hexachloroethane	8.65	117	179151	38.886	ng	98
19) n-Nitroso-di-n-propylamine	8.39	70	382045	39.398	ng	99
20) 3+4-Methylphenols	8.35	107	436842	40.297	ng	98
22) Acetophenone	8.39	105	635462	38.558	ng	99
24) Nitrobenzene	8.76	77	550521	38.591	ng	99
25) Isophorone	9.30	82	945984	39.241	ng	99
26) 2-Nitrophenol	9.47	139	215716	40.736	ng	95
27) 2,4-Dimethylphenol	9.55	122	357553	39.536	ng	96
28) bis(2-Chloroethoxy)methane	9.77	93	544344	39.359	ng	100
29) 2,4-Dichlorophenol	10.00	162	410373	39.410	ng	98
30) 1,2,4-Trichlorobenzene	10.22	180	475640	38.256	ng	98
31) Naphthalene	10.40	128	1211185	38.623	ng	100
32) Benzoic acid	9.70	122	234831	37.626	ng	95
33) 4-Chloroaniline	10.51	127	526938	39.579	ng	99
34) Hexachlorobutadiene	10.69	225	313122	37.269	ng	98
35) Caprolactam	11.30	113	122535	39.985	ng	97
36) 4-Chloro-3-methylphenol	11.64	107	442600	39.561	ng	97
37) 2-Methylnaphthalene	12.02	142	964116	39.236	ng	99
38) 1-Methylnaphthalene	12.24	142	893060	38.708	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.40	216	584195	38.453	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.38	237	399062	39.911	nq	98
43) 2,4,6-Trichlorophenol	12.64	196	370132	39.737	nq	97
44) 2,4,5-Trichlorophenol	12.72	196	404980	39.996	nq	97
46) 1,1'-Biphenyl	13.05	154	1299358	38.905	nq	100
47) 2-Chloronaphthalene	13.09	162	1030090	38.899	nq	100
48) 2-Nitroaniline	13.29	65	330046	40.759	nq	99
49) Acenaphthylene	13.94	152	1588096	39.587	nq	100
50) Dimethylphthalate	13.69	163	1342158	38.695	nq	100
51) 2,6-Dinitrotoluene	13.80	165	297700	40.292	nq	99
52) Acenaphthene	14.29	154	968999	38.934	nq	99
53) 3-Nitroaniline	14.13	138	276473	38.184	nq	99
54) 2,4-Dinitrophenol	14.34	184	167820	37.671	nq	94
55) Dibenzofuran	14.62	168	1614892	38.358	nq	99
56) 4-Nitrophenol	14.44	139	237468	40.632	nq	97
57) 2,4-Dinitrotoluene	14.59	165	423356	41.350	nq	98
58) Fluorene	15.27	166	1375929	38.937	nq	99
59) 2,3,4,6-Tetrachlorophenol	14.86	232	386060	39.440	nq	99
60) Diethylphthalate	15.06	149	1384845	38.817	nq	99
61) 4-Chlorophenyl-phenylether	15.27	204	744037	38.327	nq	100
62) 4-Nitroaniline	15.29	138	297185	38.386	nq	98
63) Azobenzene	15.57	77	1423626	38.580	nq	99
65) 4,6-Dinitro-2-methylphenol	15.36	198	249744	39.567	nq	98
66) n-Nitrosodiphenylamine	15.49	169	1182767	38.877	nq	99
67) 4-Bromophenyl-phenylether	16.17	248	465253	38.315	nq	99
68) Hexachlorobenzene	16.28	284	545620	38.494	nq	99
69) Atrazine	16.45	200	349054	39.223	nq	99
70) Pentachlorophenol	16.63	266	335184	39.104	nq	99
71) Phenanthrene	17.01	178	2215801	38.745	nq	100
72) Anthracene	17.10	178	2271076	39.348	nq	99
73) Carbazole	17.37	167	2013533	39.963	nq	100
74) Di-n-butylphthalate	17.96	149	2427386	40.050	nq	99
75) Fluoranthene	19.03	202	2766336	39.875	nq	100
77) Benzidine	19.22	184	1226912	37.144	nq	99
78) Pyrene	19.40	202	2834217	38.019	nq	99
80) Butylbenzylphthalate	20.32	149	1119504	36.345	nq	98
81) Benzo(a)anthracene	21.16	228	2857645	39.157	nq	100
82) 3,3'-Dichlorobenzidine	21.10	252	1145568	37.833	nq	99
83) Chrysene	21.21	228	2744077	38.692	nq	100
84) Bis(2-ethylhexyl)phthalate	21.11	149	1692167	41.790	nq	99
85) Di-n-octyl phthalate	21.98	149	2842945	42.181	nq	100
87) Indeno(1,2,3-cd)pyrene	25.52	276	3008035	40.370	nq	99
88) Benzo(b)fluoranthene	22.70	252	2993045	39.159	nq	100
89) Benzo(k)fluoranthene	22.74	252	2808664	38.010	nq	99
90) Benzo(a)pyrene	23.26	252	2536776	39.242	nq	100
91) Dibenzo(a,h)anthracene	25.53	278	2542031	40.389	nq	100
92) Benzo(g,h,i)perylene	26.18	276	2409303	39.723	nq	99

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

