

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031819\
 Data File : BM019216.D
 Acq On : 18 Mar 2019 15:27
 Operator : JU/SJ
 Sample : PB118000BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB118000BS

Manual Integrations
 APPROVED

Sohil
 3/19/2019 5:28:00 PM

Quant Time: Mar 19 07:31:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 11 17:30:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	233257	20.00	ng	-0.01
21) Naphthalene-d8	10.43	136	1033985	20.00	ng	-0.02
39) Acenaphthene-d10	14.30	164	601437	20.00	ng	-0.02
64) Phenanthrene-d10	17.06	188	1259536	20.00	ng	-0.01
76) Chrysene-d12	21.26	240	1031358	20.00	ng	-0.01
87) Perylene-d12	23.49	264	911770	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.25	112	2041799	141.76	ng	0.00
7) Phenol-d6	6.84	99	2886351	137.00	ng	0.00
23) Nitrobenzene-d5	8.82	82	1709807	78.17	ng	-0.01
42) 2,4,6-Tribromophenol	15.81	330	1132209	127.50	ng	-0.01
45) 2-Fluorobiphenyl	12.92	172	3619453	78.28	ng	-0.02
79) Terphenyl-d14	19.70	244	5001627	96.27	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.22	88	204792	31.962	ng	98
3) Pyridine	3.61	79	491273	28.182	ng	99
4) n-Nitrosodimethylamine	3.53	42	182606	33.696	ng	95
6) Aniline	7.00	93	876558	32.226	ng	99
8) 2-Chlorophenol	7.22	128	718049	41.486	ng	99
9) Benzaldehyde	6.81	77	363031	27.395	ng	99
10) Phenol	6.86	94	963816	42.449	ng	99
11) bis(2-Chloroethyl)ether	7.09	93	640320	36.959	ng	98
12) 1,3-Dichlorobenzene	7.53	146	656376	35.835	ng	100
13) 1,4-Dichlorobenzene	7.69	146	682621	36.554	ng	99
14) 1,2-Dichlorobenzene	8.00	146	671359	36.952	ng	99
15) Benzyl Alcohol	7.90	79	678196	38.895	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.18	45	639202	36.133	ng	99
17) 2-Methylphenol	8.10	107	619767	41.889	ng	99
18) Hexachloroethane	8.70	117	247297	35.639	ng	98
19) n-Nitroso-di-n-propylamine	8.46	70	624764	36.147	ng	99
20) 3+4-Methylphenols	8.43	107	834043	41.529	ng	99
22) Acetophenone	8.48	105	1036784	36.283	ng	# 99
24) Nitrobenzene	8.86	77	884302	38.872	ng	99
25) Isophorone	9.38	82	1557621	37.318	ng	100
26) 2-Nitrophenol	9.56	139	382730	40.587	ng	98
27) 2,4-Dimethylphenol	9.62	122	656471	46.828	ng	98
28) bis(2-Chloroethoxy)methane	9.86	93	939394	38.010	ng	99
29) 2,4-Dichlorophenol	10.09	162	673517	43.380	ng	99
30) 1,2,4-Trichlorobenzene	10.30	180	693342	39.604	ng	100
31) Naphthalene	10.49	128	2102856	38.586	ng	100
32) Benzoic acid	9.80	122	423429m	35.664	ng	
33) 4-Chloroaniline	10.62	127	454963	20.365	ng	99
34) Hexachlorobutadiene	10.75	225	371694	36.997	ng	99
35) Caprolactam	11.44	113	212272m	38.380	ng	
36) 4-Chloro-3-methylphenol	11.74	107	755604	39.442	ng	98
37) 2-Methylnaphthalene	12.10	142	1574832	39.766	ng	100
38) 1-Methylnaphthalene	12.33	142	1495540	38.313	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.47	216	748380	39.337	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.43	237	715179	82.991	ng	98
43) 2,4,6-Trichlorophenol	12.73	196	536735	43.656	ng	100
44) 2,4,5-Trichlorophenol	12.80	196	555140	42.225	ng	99
46) 1,1'-Biphenyl	13.13	154	2038092	38.803	ng	98
47) 2-Chloronaphthalene	13.17	162	1571427	40.081	ng	99
48) 2-Nitroaniline	13.40	65	519011	39.473	ng	97
49) Acenaphthylene	14.03	152	2534897	38.239	ng	99
50) Dimethylphthalate	13.77	163	1950267	37.906	ng	99
51) 2,6-Dinitrotoluene	13.90	165	433564	38.982	ng	97
52) Acenaphthene	14.37	154	1459225	38.308	ng	100
53) 3-Nitroaniline	14.24	138	307110	26.071	ng	95
54) 2,4-Dinitrophenol	14.45	184	379409	58.740	ng	98
55) Dibenzofuran	14.71	168	2360362	38.424	ng	97
56) 4-Nitrophenol	14.56	139	641661	66.717	ng	97
57) 2,4-Dinitrotoluene	14.70	165	601238	37.241	ng	97
58) Fluorene	15.36	166	1931333	38.142	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.94	232	483925	39.538	ng	100
60) Diethylphthalate	15.14	149	1981291	38.023	ng	99
61) 4-Chlorophenyl-phenylether	15.36	204	959400	39.012	ng	97
62) 4-Nitroaniline	15.42	138	384674	32.404	ng	97
63) Azobenzene	15.65	77	2100190	38.302	ng	99
65) 4,6-Dinitro-2-methylphenol	15.46	198	279188	33.826	ng	99
66) n-Nitrosodiphenylamine	15.58	169	1657024	41.323	ng	99
67) 4-Bromophenyl-phenylether	16.25	248	564995	40.678	ng	97
68) Hexachlorobenzene	16.35	284	659001	39.908	ng	98
69) Atrazine	16.54	200	526774	39.025	ng	99
70) Pentachlorophenol	16.71	266	740524	74.433	ng	99
71) Phenanthrene	17.10	178	2855133	38.674	ng	100
72) Anthracene	17.19	178	2913103	40.307	ng	100
73) Carbazole	17.47	167	2538275	37.211	ng	99
74) Di-n-butylphthalate	18.03	149	3327584	40.341	ng	100
75) Fluoranthene	19.13	202	3040434	35.891	ng	99
77) Benzidine	19.33	184	1381258	47.223	ng	99
78) Pyrene	19.49	202	3057752	46.858	ng	99
80) Butylbenzylphthalate	20.40	149	1471014	51.756	ng	100
81) Benzo(a)anthracene	21.24	228	2509975	38.792	ng	100
82) 3,3'-Dichlorobenzidine	21.19	252	729802	29.279	ng	99
83) Chrysene	21.30	228	2358903	37.684	ng	100
84) Bis(2-ethylhexyl)phthalate	21.17	149	2147697	51.940	ng	# 98
85) Di-n-octyl phthalate	22.04	149	3366796	48.354	ng	100
86) Indeno(1,2,3-cd)pyrene	25.76	276	2682826	39.814	ng	100
88) Benzo(b)fluoranthene	22.82	252	2328432	39.411	ng	100
89) Benzo(k)fluoranthene	22.87	252	2238902	41.201	ng	99
90) Benzo(a)pyrene	23.40	252	2194641	41.335	ng	100
91) Dibenzo(a,h)anthracene	25.76	278	2299893	45.283	ng	100
92) Benzo(g,h,i)perylene	26.45	276	2192199	47.013	ng	99

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

