

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM082019\
 Data File : BM022197.D
 Acq On : 19 Aug 2019 19:40
 Operator : HP/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

mohammad
 8/22/2019 12:48:25 PM

Quant Time: Aug 20 01:52:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM082019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 20 01:26:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	35969	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	154094	20.00	ng	0.00
39) Acenaphthene-d10	14.23	164	97369	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	188066	20.00	ng	0.00
76) Chrysene-d12	21.20	240	191220	20.00	ng	0.00
87) Perylene-d12	23.40	264	189115	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	193772	97.55	ng	0.00
7) Phenol-d6	6.76	99	262756	90.83	ng	0.00
23) Nitrobenzene-d5	8.75	82	310063	104.87	ng	0.00
42) 2,4,6-Tribromophenol	15.74	330	146180	80.01	ng	0.00
45) 2-Fluorobiphenyl	12.84	172	743059	93.79	ng	0.00
79) Terphenyl-d14	19.63	244	1234893	120.62	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.17	88	40354	53.921	ng	# 89
3) Pyridine	3.56	79	109940	51.085	ng	# 88
4) n-Nitrosodimethylamine	3.49	42	63134	76.188	ng	# 62
6) Aniline	6.93	93	176030	45.689	ng	94
8) 2-Chlorophenol	7.14	128	116696	47.733	ng	96
9) Benzaldehyde	6.75	77	80377	52.524	ng	90
10) Phenol	6.79	94	138025	46.705	ng	83
11) bis(2-Chloroethyl)ether	7.02	93	109168	46.658	ng	97
12) 1,3-Dichlorobenzene	7.46	146	140551	47.453	ng	# 93
13) 1,4-Dichlorobenzene	7.60	146	141899	47.121	ng	98
14) 1,2-Dichlorobenzene	7.92	146	139974	47.431	ng	96
15) Benzyl Alcohol	7.83	79	114377m	48.833	ng	
16) 2,2'-oxybis(1-Chloropropan	8.10	45	152343	50.338	ng	98
17) 2-Methylphenol	8.02	107	99760	47.258	ng	# 93
18) Hexachloroethane	8.62	117	62427	57.738	ng	90
19) n-Nitroso-di-n-propylamine	8.39	70	103589	49.208	ng	91
20) 3+4-Methylphenols	8.36	107	134572	46.313	ng	94
22) Acetophenone	8.41	105	189385	48.229	ng	# 90
24) Nitrobenzene	8.79	77	159819	54.567	ng	93
25) Isophorone	9.30	82	270042	48.812	ng	95
26) 2-Nitrophenol	9.49	139	68104	47.488	ng	# 73
27) 2,4-Dimethylphenol	9.55	122	101367	48.461	ng	93
28) bis(2-Chloroethoxy)methane	9.78	93	162994	47.414	ng	99
29) 2,4-Dichlorophenol	10.01	162	122710	45.328	ng	96
30) 1,2,4-Trichlorobenzene	10.21	180	148901	46.660	ng	99
31) Naphthalene	10.40	128	384166	47.110	ng	100
32) Benzoic acid	9.75	122	90135m	53.266	ng	
33) 4-Chloroaniline	10.53	127	162932	44.013	ng	# 95
34) Hexachlorobutadiene	10.65	225	126735	62.898	ng	98
35) Caprolactam	11.37	113	32615m	38.518	ng	
36) 4-Chloro-3-methylphenol	11.66	107	126460	45.907	ng	97
37) 2-Methylnaphthalene	12.02	142	279087	44.479	ng	# 93
38) 1-Methylnaphthalene	12.24	142	268703	44.898	ng	# 99
40) 1,2,4,5-Tetrachlorobenzene	12.39	216	185644	50.372	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.34	237	92829	43.450	ng	97
43) 2,4,6-Trichlorophenol	12.65	196	112105	45.969	ng	96
44) 2,4,5-Trichlorophenol	12.73	196	113326	44.874	ng #	93
46) 1,1'-Biphenyl	13.05	154	370306	44.435	ng	99
47) 2-Chloronaphthalene	13.10	162	303771	44.880	ng	96
48) 2-Nitroaniline	13.33	65	97800	52.818	ng	84
49) Acenaphthylene	13.95	152	463138	44.964	ng	99
50) Dimethylphthalate	13.70	163	391595	45.234	ng	100
51) 2,6-Dinitrotoluene	13.84	165	77024	41.357	ng #	63
52) Acenaphthene	14.29	154	270217	43.482	ng	96
53) 3-Nitroaniline	14.18	138	78435	41.310	ng	95
54) 2,4-Dinitrophenol	14.40	184	41602	41.584	ng #	65
55) Dibenzofuran	14.63	168	443885	43.785	ng #	91
56) 4-Nitrophenol	14.50	139	52053	35.852	ng #	58
57) 2,4-Dinitrotoluene	14.64	165	109913	42.740	ng #	65
58) Fluorene	15.29	166	369930	44.660	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.87	232	112326	46.552	ng #	94
60) Diethylphthalate	15.07	149	407909	47.841	ng	96
61) 4-Chlorophenyl-phenylether	15.29	204	234454	50.102	ng	96
62) 4-Nitroaniline	15.36	138	72553	36.408	ng #	59
63) Azobenzene	15.58	77	397230	56.797	ng	84
65) 4,6-Dinitro-2-methylphenol	15.41	198	56925	49.091	ng	83
66) n-Nitrosodiphenylamine	15.51	169	295282	49.813	ng	98
67) 4-Bromophenyl-phenylether	16.18	248	136437	54.497	ng #	92
68) Hexachlorobenzene	16.29	284	153818	51.429	ng #	84
69) Atrazine	16.47	200	106702	55.354	ng	98
70) Pentachlorophenol	16.64	266	78100	48.465	ng	98
71) Phenanthrene	17.03	178	528085	48.723	ng	99
72) Anthracene	17.13	178	525612	48.956	ng	99
73) Carbazole	17.42	167	435179	45.842	ng #	97
74) Di-n-butylphthalate	17.96	149	661538	56.925	ng #	98
75) Fluoranthene	19.06	202	628341	50.202	ng	99
77) Benzidine	19.27	184	201298	36.252	ng	99
78) Pyrene	19.43	202	631270	45.520	ng	100
80) Butylbenzylphthalate	20.33	149	294454	53.270	ng	86
81) Benzo(a)anthracene	21.18	228	638933	49.016	ng	98
82) 3,3'-Dichlorobenzidine	21.13	252	227592	46.098	ng	100
83) Chrysene	21.24	228	619885	49.920	ng	99
84) Bis(2-ethylhexyl)phthalate	21.10	149	443454	56.368	ng	95
85) Di-n-octyl phthalate	21.97	149	721192	58.290	ng #	94
86) Indeno(1,2,3-cd)pyrene	25.60	276	712741	51.872	ng	97
88) Benzo(b)fluoranthene	22.74	252	622397	48.784	ng	99
89) Benzo(k)fluoranthene	22.79	252	591946	47.998	ng	99
90) Benzo(a)pyrene	23.30	252	558177	48.486	ng	99
91) Dibenzo(a,h)anthracene	25.60	278	593604	51.719	ng	97
92) Benzo(g,h,i)perylene	26.27	276	519298	48.518	ng	97

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

