

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122320\
 Data File : BM028244.D
 Acq On : 23 Dec 2020 09:50
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 12/24/2020 10:43:30 AM

Quant Time: Dec 23 13:05:50 2020
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM121620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 23 12:17:02 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.198	152	159369	20.00	ng	0.00
21) Naphthalene-d8	11.028	136	649793	20.00	ng	# 0.00
39) Acenaphthene-d10	14.827	164	385416	20.00	ng	0.00
64) Phenanthrene-d10	17.545	188	787242	20.00	ng	# 0.00
76) Chrysene-d12	21.662	240	629678	20.00	ng	# 0.00
86) Perylene-d12	24.021	264	531127	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.693	112	694845	71.85	ng	0.00
7) Phenol-d6	7.340	99	986985	71.47	ng	0.00
23) Nitrobenzene-d5	9.375	82	927633	70.84	ng	0.00
42) 2,4,6-Tribromophenol	16.304	330	377288	73.74	ng	0.00
45) 2-Fluorobiphenyl	13.457	172	2110548	71.19	ng	0.00
79) Terphenyl-d14	20.121	244	2818039	81.21	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.446	88	135813	34.34	ng	# 70
3) Pyridine	3.887	79	392125	34.57	ng	# 52
4) n-Nitrosodimethylamine	3.787	42	199835	32.98	ng	# 62
6) Aniline	7.510	93	541614	34.92	ng	98
8) 2-Chlorophenol	7.751	128	399994	35.95	ng	97
9) Benzaldehyde	7.316	77	253106	32.80	ng	82
10) Phenol	7.369	94	462709	35.27	ng	85
11) bis(2-Chloroethyl)ether	7.604	93	382272	34.89	ng	92
12) 1,3-Dichlorobenzene	8.081	146	469193	35.88	ng	# 92
13) 1,4-Dichlorobenzene	8.234	146	474533	35.83	ng	98
14) 1,2-Dichlorobenzene	8.557	146	458786	35.94	ng	98
15) Benzyl Alcohol	8.439	79	340323	34.77	ng	# 85
16) 2,2'-oxybis(1-Chloropr...	8.734	45	665028	33.60	ng	70
17) 2-Methylphenol	8.634	107	330177	35.98	ng	# 88
18) Hexachloroethane	9.292	117	177487	35.99	ng	91
19) n-Nitroso-di-n-propyla...	9.016	70	303108	34.81	ng	# 68
20) 3+4-Methylphenols	8.969	107	442245	35.68	ng	90
22) Acetophenone	9.034	105	580774	34.79	ng	# 89
24) Nitrobenzene	9.416	77	449223	35.17	ng	# 88
25) Isophorone	9.945	82	766231	35.53	ng	96
26) 2-Nitrophenol	10.133	139	213890	38.21	ng	# 80
27) 2,4-Dimethylphenol	10.180	122	315142	35.50	ng	88
28) bis(2-Chloroethoxy)met...	10.422	93	532301	34.38	ng	99
29) 2,4-Dichlorophenol	10.663	162	373476	36.17	ng	97
30) 1,2,4-Trichlorobenzene	10.886	180	425676	35.19	ng	98
31) Naphthalene	11.075	128	1295301	35.39	ng	99
32) Benzoic acid	10.328	122	270914	38.10	ng	# 82
33) 4-Chloroaniline	11.180	127	545239	34.89	ng	# 91
34) Hexachlorobutadiene	11.351	225	254867	35.50	ng	98
35) Caprolactam	11.974	113	121279	36.15	ng	# 55
36) 4-Chloro-3-methylphenol	12.286	107	391378	35.42	ng	93
37) 2-Methylnaphthalene	12.669	142	907674	35.04	ng	96
38) 1-Methylnaphthalene	12.886	142	860741	34.91	ng	100
40) 1,2,4,5-Tetrachloroben...	13.033	216	437359	35.78	ng	99
41) Hexachlorocyclopentadiene	13.010	237	219960	34.88	ng	99
43) 2,4,6-Trichlorophenol	13.269	196	300747	36.64	ng	97

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44) 2,4,5-Trichlorophenol	13.339	196	310591	36.08	ng	96
46) 1,1'-Biphenyl	13.669	154	1135751	35.29	ng	99
47) 2-Chloronaphthalene	13.716	162	922809	35.40	ng	99
48) 2-Nitroaniline	13.910	65	290944	36.74	ng	88
49) Acenaphthylene	14.551	152	1401248	35.88	ng	100
50) Dimethylphthalate	14.274	163	1105775	34.57	ng	100
51) 2,6-Dinitrotoluene	14.398	165	239221	35.77	ng	90
52) Acenaphthene	14.892	154	933019m	35.05	ng	
53) 3-Nitroaniline	14.727	138	272773	36.06	ng	87
54) 2,4-Dinitrophenol	14.927	184	132250	34.60	ng	91
55) Dibenzofuran	15.221	168	1327828	35.03	ng	97
56) 4-Nitrophenol	15.015	139	210838	35.25	ng	# 73
57) 2,4-Dinitrotoluene	15.180	165	322272	36.39	ng	91
58) Fluorene	15.862	166	1088020	35.03	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.439	232	282389	35.74	ng	94
60) Diethylphthalate	15.627	149	1146448	34.82	ng	99
61) 4-Chlorophenyl-phenyle...	15.851	204	539581	34.78	ng	96
62) 4-Nitroaniline	15.880	138	293088	35.73	ng	# 78
63) Azobenzene	16.139	77	1068005	35.18	ng	91
65) 4,6-Dinitro-2-methylph...	15.939	198	166888	35.90	ng	85
66) n-Nitrosodiphenylamine	16.062	169	938050	35.87	ng	98
67) 4-Bromophenyl-phenylether	16.739	248	336341	35.73	ng	97
68) Hexachlorobenzene	16.857	284	384554	35.70	ng	96
69) Atrazine	16.998	200	308027	36.40	ng	99
70) Pentachlorophenol	17.192	266	225832	37.49	ng	98
71) Phenanthrene	17.592	178	1671060	34.95	ng	100
72) Anthracene	17.680	178	1638920	35.46	ng	99
73) Carbazole	17.945	167	1516913	34.79	ng	99
74) Di-n-butylphthalate	18.480	149	2000828	36.71	ng	98
75) Fluoranthene	19.574	202	1807269	33.76	ng	96
77) Benzidine	19.745	184	693223	43.64	ng	99
78) Pyrene	19.933	202	1824060	40.01	ng	99
80) Butylbenzylphthalate	20.792	149	848712	42.75	ng	98
81) Benzo(a)anthracene	21.644	228	1520985	35.46	ng	100
82) 3,3'-Dichlorobenzidine	21.568	252	510957	36.93	ng	# 99
83) Chrysene	21.697	228	1439432	34.66	ng	100
84) Bis(2-ethylhexyl)phtha...	21.550	149	1212413	42.67	ng	98
85) Di-n-octyl phthalate	22.462	149	1874348	40.70	ng	# 91
87) Indeno(1,2,3-cd)pyrene	26.456	276	1308276	37.05	ng	# 89
88) Benzo(b)fluoranthene	23.309	252	1320428	36.49	ng	# 96
89) Benzo(k)fluoranthene	23.356	252	1230099	34.98	ng	# 97
90) Benzo(a)pyrene	23.921	252	1149228	35.39	ng	# 96
91) Dibenzo(a,h)anthracene	26.462	278	1083013	36.76	ng	# 94
92) Benzo(g,h,i)perylene	27.203	276	1059396	37.64	ng	# 91

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

