

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM102318\
 Data File : BM017285.D
 Acq On : 23 Oct 2018 15:22
 Operator : MJ/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 10/24/2018 10:57:24 AM

Quant Time: Oct 24 07:02:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM101718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 17 14:23:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	250467	20.00	ng	-0.01
21) Naphthalene-d8	10.43	136	1119449	20.00	ng	-0.02
39) Acenaphthene-d10	14.30	164	656411	20.00	ng	-0.01
64) Phenanthrene-d10	17.06	188	1452457	20.00	ng	-0.01
76) Chrysene-d12	21.25	240	1660279	20.00	ng	-0.01
87) Perylene-d12	23.46	264	1506330	20.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.28	112	1329979	89.81	ng	0.00
7) Phenol-d6	6.85	99	1799441	89.22	ng	0.00
23) Nitrobenzene-d5	8.82	82	1736096	90.70	ng	-0.02
42) 2,4,6-Tribromophenol	15.80	330	830174	93.52	ng	-0.01
45) 2-Fluorobiphenyl	12.92	172	4136210	83.84	ng	-0.01
79) Terphenyl-d14	19.69	244	6690388	90.83	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.26	88	302202	39.963	ng	# 86
3) Pyridine	3.65	79	761353	43.757	ng	# 87
4) n-Nitrosodimethylamine	3.56	42	288900	41.361	ng	# 78
6) Aniline	7.00	93	1055003	41.820	ng	98
8) 2-Chlorophenol	7.23	128	772315	45.080	ng	98
9) Benzaldehyde	6.82	77	494976	38.490	ng	97
10) Phenol	6.87	94	945504	43.559	ng	99
11) bis(2-Chloroethyl)ether	7.10	93	753894	43.568	ng	99
12) 1,3-Dichlorobenzene	7.55	146	856841	42.904	ng	99
13) 1,4-Dichlorobenzene	7.70	146	868370	42.558	ng	98
14) 1,2-Dichlorobenzene	8.01	146	847635	43.117	ng	99
15) Benzyl Alcohol	7.91	79	716556	48.605	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.19	45	993601	41.101	ng	98
17) 2-Methylphenol	8.11	107	635627	45.622	ng	99
18) Hexachloroethane	8.72	117	309357	44.303	ng	99
19) n-Nitroso-di-n-propylamine	8.47	70	619899	46.678	ng	98
20) 3+4-Methylphenols	8.44	107	896435	47.042	ng	98
22) Acetophenone	8.48	105	1195438	42.128	ng	# 100
24) Nitrobenzene	8.86	77	878386	43.385	ng	94
25) Isophorone	9.38	82	1452827	45.966	ng	98
26) 2-Nitrophenol	9.56	139	462175	48.683	ng	96
27) 2,4-Dimethylphenol	9.63	122	642857	45.998	ng	100
28) bis(2-Chloroethoxy)methane	9.86	93	972193	43.702	ng	99
29) 2,4-Dichlorophenol	10.09	162	762058	46.963	ng	99
30) 1,2,4-Trichlorobenzene	10.30	180	832919	42.788	ng	98
31) Naphthalene	10.49	128	2286014	41.670	ng	100
32) Benzoic acid	9.80	122	628696	60.290	ng	93
33) 4-Chloroaniline	10.60	127	987373	41.673	ng	98
34) Hexachlorobutadiene	10.76	225	512884	41.583	ng	97
35) Caprolactam	11.43	113	298459	50.469	ng	90
36) 4-Chloro-3-methylphenol	11.74	107	857046	46.856	ng	96
37) 2-Methylnaphthalene	12.10	142	1639182	43.664	ng	99
38) 1-Methylnaphthalene	12.33	142	1585454	43.392	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.47	216	982884	42.519	ng	100

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41) Hexachlorocyclopentadiene	12.45	237	527804	47.229	ng	100
43) 2,4,6-Trichlorophenol	12.73	196	649208	48.379	ng	99
44) 2,4,5-Trichlorophenol	12.80	196	684877	47.594	ng	97
46) 1,1'-Biphenyl	13.13	154	2475286	41.904	ng	100
47) 2-Chloronaphthalene	13.17	162	1832535	42.170	ng	99
48) 2-Nitroaniline	13.39	65	581253	47.422	ng	98
49) Acenaphthylene	14.02	152	2751729	43.685	ng	99
50) Dimethylphthalate	13.77	163	2292232	45.263	ng	100
51) 2,6-Dinitrotoluene	13.89	165	518378	46.507	ng	99
52) Acenaphthene	14.37	154	1675274	42.359	ng	97
53) 3-Nitroaniline	14.23	138	485399	40.209	ng	98
54) 2,4-Dinitrophenol	14.43	184	316433	51.642	ng	96
55) Dibenzofuran	14.70	168	2712410	41.241	ng	99
56) 4-Nitrophenol	14.56	139	476996	45.515	ng	99
57) 2,4-Dinitrotoluene	14.68	165	731127	48.804	ng	# 98
58) Fluorene	15.36	166	2072516	42.573	ng	98
59) 2,3,4,6-Tetrachlorophenol	14.94	232	610351	46.532	ng	98
60) Diethylphthalate	15.14	149	2922308	58.187	ng	100
61) 4-Chlorophenyl-phenylether	15.36	204	1161743	41.860	ng	96
62) 4-Nitroaniline	15.39	138	519827	39.518	ng	98
63) Azobenzene	15.65	77	2131185	43.658	ng	99
65) 4,6-Dinitro-2-methylphenol	15.44	198	457401	48.889	ng	97
66) n-Nitrosodiphenylamine	15.57	169	2046759	46.577	ng	99
67) 4-Bromophenyl-phenylether	16.25	248	752131	47.161	ng	96
68) Hexachlorobenzene	16.36	284	829235	44.762	ng	97
69) Atrazine	16.54	200	749277	47.641	ng	98
70) Pentachlorophenol	16.71	266	627935	49.481	ng	99
71) Phenanthrene	17.10	178	3296584	42.067	ng	99
72) Anthracene	17.19	178	3390093	45.529	ng	99
73) Carbazole	17.47	167	3471491	45.641	ng	99
74) Di-n-butylphthalate	18.03	149	4255819	53.049	ng	99
75) Fluoranthene	19.12	202	4038194	45.794	ng	99
78) Pyrene	19.49	202	4206989	45.337	ng	100
80) Butylbenzylphthalate	20.40	149	2021841	55.874	ng	99
81) Benzo(a)anthracene	21.23	228	4115083	44.044	ng	100
82) 3,3'-Dichlorobenzidine	21.17	252	1140920	32.764	ng	98
83) Chrysene	21.29	228	3894590	42.114	ng	99
84) Bis(2-ethylhexyl)phthalate	21.17	149	2929257	53.532	ng	98
85) Di-n-octyl phthalate	22.05	149	4904552	52.938	ng	97
86) Indeno(1,2,3-cd)pyrene	25.67	276	3538256	34.208	ng	99
88) Benzo(b)fluoranthene	22.80	252	3585774m	43.540	ng	
89) Benzo(k)fluoranthene	22.84	252	3652858	44.203	ng	100
90) Benzo(a)pyrene	23.37	252	3364954	43.033	ng	99
91) Dibenzo(a,h)anthracene	25.69	278	3015108	38.832	ng	99
92) Benzo(g,h,i)perylene	26.35	276	2872554	37.328	ng	99

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

