

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102918\
 Data File : BM017408.D
 Acq On : 29 Oct 2018 15:04
 Operator : MJ/SJ
 Sample : J5651-03MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RBR-200030MSD

Manual Integrations
 APPROVED

Sohil

10/30/2018 4:43:05 PM

Quant Time: Oct 30 03:07:40 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM101718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 29 15:08:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	288774	20.00	ng	0.00
21) Naphthalene-d8	10.42	136	1271667	20.00	ng	0.00
39) Acenaphthene-d10	14.29	164	789042	20.00	ng	0.00
64) Phenanthrene-d10	17.04	188	1895468	20.00	ng	0.00
76) Chrysene-d12	21.24	240	2246506	20.00	ng	0.00
87) Perylene-d12	23.45	264	2037845	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.27	112	1856379	108.73	ng	0.00
7) Phenol-d6	6.84	99	2403168	103.35	ng	0.00
23) Nitrobenzene-d5	8.80	82	1684776	77.49	ng	0.00
42) 2,4,6-Tribromophenol	15.79	330	1371361	128.52	ng	0.00
45) 2-Fluorobiphenyl	12.90	172	4253501	71.73	ng	0.00
79) Terphenyl-d14	19.69	244	7335426	73.60	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.26	88	217152	24.907	ng	# 80
3) Pyridine	3.65	79	605224	30.170	ng	88
4) n-Nitrosodimethylamine	3.56	42	331687	41.187	ng	# 74
6) Aniline	6.99	93	437494	15.042	ng	97
8) 2-Chlorophenol	7.22	128	936305	47.402	ng	99
9) Benzaldehyde	6.80	77	160281	10.810	ng	96
10) Phenol	6.86	94	1030004	41.157	ng	98
11) bis(2-Chloroethyl)ether	7.09	93	822842	41.245	ng	99
12) 1,3-Dichlorobenzene	7.53	146	936494	40.672	ng	99
13) 1,4-Dichlorobenzene	7.68	146	947433	40.274	ng	98
14) 1,2-Dichlorobenzene	7.99	146	927541	40.923	ng	99
15) Benzyl Alcohol	7.90	79	852250	50.141	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.18	45	1088139	39.040	ng	99
17) 2-Methylphenol	8.10	107	779612	48.533	ng	98
18) Hexachloroethane	8.71	117	337354	41.903	ng	100
19) n-Nitroso-di-n-propylamine	8.46	70	704951	46.041	ng	97
20) 3+4-Methylphenols	8.42	107	1078892	49.107	ng	99
22) Acetophenone	8.47	105	1407033	43.649	ng	# 99
24) Nitrobenzene	8.85	77	1062681	46.205	ng	96
25) Isophorone	9.37	82	1957380	54.517	ng	99
26) 2-Nitrophenol	9.55	139	530373	49.122	ng	99
27) 2,4-Dimethylphenol	9.62	122	923795	58.187	ng	98
28) bis(2-Chloroethoxy)methane	9.85	93	1206194	47.730	ng	99
29) 2,4-Dichlorophenol	10.07	162	959831	52.070	ng	99
30) 1,2,4-Trichlorobenzene	10.29	180	935591	42.310	ng	99
31) Naphthalene	10.47	128	2915571	46.785	ng	99
32) Benzoic acid	9.79	122	616803	52.069	ng	97
33) 4-Chloroaniline	10.60	127	135807	5.046	ng	97
34) Hexachlorobutadiene	10.75	225	552993	39.468	ng	98
35) Caprolactam	11.40	113	272836m	40.614	ng	
36) 4-Chloro-3-methylphenol	11.73	107	1097594	52.824	ng	99
37) 2-Methylnaphthalene	12.09	142	2243807	52.616	ng	99
38) 1-Methylnaphthalene	12.31	142	2138201	51.515	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.46	216	1137163	40.924	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.44	237	1305966	97.218	ng	100
43) 2,4,6-Trichlorophenol	12.71	196	820152	50.844	ng	99
44) 2,4,5-Trichlorophenol	12.79	196	904422	52.287	ng	99
46) 1,1'-Biphenyl	13.12	154	2918535	41.103	ng	99
47) 2-Chloronaphthalene	13.16	162	2268040	43.419	ng	99
48) 2-Nitroaniline	13.37	65	695318	47.220	ng	96
49) Acenaphthylene	14.01	152	3822815	50.488	ng	100
50) Dimethylphthalate	13.76	163	3020422	49.617	ng	99
51) 2,6-Dinitrotoluene	13.89	165	684438	51.084	ng	96
52) Acenaphthene	14.36	154	2247266	47.270	ng	97
53) 3-Nitroaniline	14.22	138	398987	27.495	ng	98
54) 2,4-Dinitrophenol	14.43	184	932839	114.702	ng	96
55) Dibenzofuran	14.70	168	3634357	45.971	ng	99
56) 4-Nitrophenol	14.54	139	1633990	120.983	ng	98
57) 2,4-Dinitrotoluene	14.67	165	1000478	55.558	ng	# 98
58) Fluorene	15.34	166	2985153	51.013	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.93	232	858895	54.474	ng	99
60) Diethylphthalate	15.14	149	3092215	51.220	ng	99
61) 4-Chlorophenyl-phenylether	15.34	204	1489226	44.640	ng	96
62) 4-Nitroaniline	15.39	138	721310	44.884	ng	99
63) Azobenzene	15.64	77	2852809	48.617	ng	99
65) 4,6-Dinitro-2-methylphenol	15.44	198	615049	50.161	ng	98
66) n-Nitrosodiphenylamine	15.56	169	2709757	47.252	ng	100
67) 4-Bromophenyl-phenylether	16.24	248	997047	47.907	ng	96
68) Hexachlorobenzene	16.34	284	1076608	44.533	ng	99
69) Atrazine	16.53	200	1009082	49.164	ng	98
70) Pentachlorophenol	16.70	266	1419652	85.722	ng	99
71) Phenanthrene	17.09	178	4915761	48.068	ng	100
72) Anthracene	17.18	178	5059683	52.069	ng	100
73) Carbazole	17.46	167	4763266	47.987	ng	100
74) Di-n-butylphthalate	18.02	149	5547293	52.987	ng	99
75) Fluoranthene	19.11	202	6046023	52.539	ng	98
77) Benzidine	19.31	184	388239	6.816	ng	98
78) Pyrene	19.47	202	6129872	48.821	ng	99
80) Butylbenzylphthalate	20.39	149	2758328	56.276	ng	95
81) Benzo(a)anthracene	21.23	228	6379344	50.461	ng	99
82) 3,3'-Dichlorobenzidine	21.17	252	1070120	22.712	ng	98
83) Chrysene	21.28	228	6004838	47.989	ng	99
84) Bis(2-ethylhexyl)phthalate	21.17	149	3935210	53.187	ng	97
85) Di-n-octyl phthalate	22.04	149	6929381	55.030	ng	99
86) Indeno(1,2,3-cd)pyrene	25.67	276	5426796	38.775	ng	99
88) Benzo(b)fluoranthene	22.79	252	6162652	55.312	ng	99
89) Benzo(k)fluoranthene	22.84	252	5772907	51.637	ng	99
90) Benzo(a)pyrene	23.36	252	5511171	52.097	ng	99
91) Dibenzo(a,h)anthracene	25.67	278	4665241	44.413	ng	99
92) Benzo(g,h,i)perylene	26.33	276	4350516	41.788	ng	99

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

