

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052419\
 Data File : BM020567.D
 Acq On : 25 May 2019 06:56
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

mohammad
 5/28/2019 8:41:37 AM

Quant Time: May 27 03:57:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM052419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 24 16:07:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	107372	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	420474	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	285107	20.00	ng	0.00
64) Phenanthrene-d10	17.07	188	675600	20.00	ng	0.00
76) Chrysene-d12	21.27	240	537056	20.00	ng	0.00
87) Perylene-d12	23.51	264	591004	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.25	112	429557	78.49	ng	0.00
7) Phenol-d6	6.85	99	645963	78.20	ng	0.00
23) Nitrobenzene-d5	8.84	82	788659	78.73	ng	0.00
42) 2,4,6-Tribromophenol	15.83	330	411237	72.00	ng	0.00
45) 2-Fluorobiphenyl	12.93	172	1874475	82.50	ng	0.00
79) Terphenyl-d14	19.71	244	2897533	97.53	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.17	88	100590	35.122	ng	98
3) Pyridine	3.58	79	303188	36.056	ng	98
4) n-Nitrosodimethylamine	3.50	42	73109	35.995	ng	# 94
6) Aniline	7.00	93	453034	37.912	ng	98
8) 2-Chlorophenol	7.23	128	263541	39.378	ng	99
9) Benzaldehyde	6.82	77	211112	40.030	ng	97
10) Phenol	6.87	94	344612	38.632	ng	99
11) bis(2-Chloroethyl)ether	7.10	93	253695	38.110	ng	99
12) 1,3-Dichlorobenzene	7.55	146	343202	40.043	ng	99
13) 1,4-Dichlorobenzene	7.70	146	340460	39.741	ng	98
14) 1,2-Dichlorobenzene	8.01	146	340186	39.609	ng	98
15) Benzyl Alcohol	7.92	79	279935m	36.621	ng	
16) 2,2'-oxybis(1-Chloropropan	8.19	45	180573	38.919	ng	95
17) 2-Methylphenol	8.11	107	225681	38.322	ng	99
18) Hexachloroethane	8.72	117	100959	29.092	ng	95
19) n-Nitroso-di-n-propylamine	8.47	70	267233	36.432	ng	99
20) 3+4-Methylphenols	8.45	107	299535	37.927	ng	98
22) Acetophenone	8.50	105	442236	38.194	ng	99
24) Nitrobenzene	8.88	77	380046	38.414	ng	99
25) Isophorone	9.40	82	700230	37.565	ng	100
26) 2-Nitrophenol	9.58	139	116732	31.977	ng	97
27) 2,4-Dimethylphenol	9.64	122	208482	38.338	ng	98
28) bis(2-Chloroethoxy)methane	9.88	93	377133	38.866	ng	99
29) 2,4-Dichlorophenol	10.11	162	291336	40.281	ng	98
30) 1,2,4-Trichlorobenzene	10.32	180	402732	40.724	ng	97
31) Naphthalene	10.50	128	865011	39.843	ng	99
32) Benzoic acid	9.83	122	122723	36.804	ng	91
33) 4-Chloroaniline	10.63	127	333262	38.459	ng	98
34) Hexachlorobutadiene	10.76	225	302797	41.183	ng	98
35) Caprolactam	11.47	113	79936	33.581	ng	93
36) 4-Chloro-3-methylphenol	11.76	107	301456	38.301	ng	96
37) 2-Methylnaphthalene	12.12	142	644866	38.880	ng	98
38) 1-Methylnaphthalene	12.35	142	620116	38.588	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.49	216	508109	42.395	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.45	237	19127	11.241	nq	97
43) 2,4,6-Trichlorophenol	12.75	196	309171	41.684	nq	99
44) 2,4,5-Trichlorophenol	12.82	196	298160	43.048	nq	99
46) 1,1'-Biphenyl	13.15	154	896911	40.847	nq	100
47) 2-Chloronaphthalene	13.19	162	768316	41.276	nq	98
48) 2-Nitroaniline	13.42	65	238570	41.706	nq	98
49) Acenaphthylene	14.05	152	1126946	39.811	nq	99
50) Dimethylphthalate	13.79	163	955221	37.825	nq	100
51) 2,6-Dinitrotoluene	13.93	165	193852	36.830	nq	100
52) Acenaphthene	14.39	154	682416	40.004	nq	99
53) 3-Nitroaniline	14.26	138	181155	39.718	nq	98
54) 2,4-Dinitrophenol	14.49	184	6631	12.594	nq	# 76
55) Dibenzofuran	14.73	168	1145816	38.816	nq	100
56) 4-Nitrophenol	14.59	139	105224	35.817	nq	94
57) 2,4-Dinitrotoluene	14.72	165	261878	34.455	nq	95
58) Fluorene	15.37	166	944846	38.532	nq	100
59) 2,3,4,6-Tetrachlorophenol	14.96	232	300956	37.343	nq	98
60) Diethylphthalate	15.15	149	936729	37.648	nq	100
61) 4-Chlorophenyl-phenylether	15.37	204	611765	39.079	nq	99
62) 4-Nitroaniline	15.43	138	164372	37.152	nq	99
63) Azobenzene	15.66	77	918253	37.927	nq	99
65) 4,6-Dinitro-2-methylphenol	15.49	198	17016	9.786	nq	96
66) n-Nitrosodiphenylamine	15.59	169	793399	43.311	nq	97
67) 4-Bromophenyl-phenylether	16.27	248	392940	43.660	nq	99
68) Hexachlorobenzene	16.37	284	460387	42.190	nq	98
69) Atrazine	16.55	200	298970	39.667	nq	99
70) Pentachlorophenol	16.73	266	212547	37.709	nq	99
71) Phenanthrene	17.12	178	1448624	39.470	nq	100
72) Anthracene	17.21	178	1389726	39.045	nq	99
73) Carbazole	17.49	167	1148934	37.614	nq	100
74) Di-n-butylphthalate	18.04	149	1567353	41.713	nq	100
75) Fluoranthene	19.14	202	1621240	31.765	nq	100
77) Benzidine	19.34	184	650258	65.884	nq	99
78) Pyrene	19.50	202	1585804	47.033	nq	100
80) Butylbenzylphthalate	20.40	149	564824	47.678	nq	97
81) Benzo(a)anthracene	21.25	228	1463657	39.377	nq	100
82) 3,3'-Dichlorobenzidine	21.19	252	594524	43.822	nq	98
83) Chrysene	21.31	228	1429303	39.795	nq	100
84) Bis(2-ethylhexyl)phthalate	21.17	149	797329	46.178	nq	98
85) Di-n-octyl phthalate	22.04	149	1259075	41.946	nq	100
86) Indeno(1,2,3-cd)pyrene	25.79	276	1712277	39.429	nq	99
88) Benzo(b)fluoranthene	22.83	252	1476827	38.179	nq	100
89) Benzo(k)fluoranthene	22.88	252	1451956	38.675	nq	99
90) Benzo(a)pyrene	23.42	252	1383769	38.874	nq	99
91) Dibenzo(a,h)anthracene	25.80	278	1447430	39.724	nq	98
92) Benzo(g,h,i)perylene	26.49	276	1359501	40.250	nq	99

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

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