

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072518\
 Data File : BN002165.D
 Acq On : 26 Jul 2018 08:34
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

Sohil
 7/27/2018 11:36:06 AM

Quant Time: Jul 26 09:12:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN072518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 25 20:08:02 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	285299	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	1209880	20.00	ng	0.00
38) Acenaphthene-d10	14.32	164	640210	20.00	ng	0.00
63) Phenanthrene-d10	17.07	188	1252386	20.00	ng	0.00
75) Chrysene-d12	21.27	240	1113071	20.00	ng	0.00
86) Perylene-d12	23.49	264	1099412	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	1371373	77.39	ng	0.00
7) Phenol-d6	6.88	99	1910501	76.82	ng	0.00
23) Nitrobenzene-d5	8.85	82	1907031	80.83	ng	0.00
41) 2,4,6-Tribromophenol	15.82	330	592524	71.27	ng	0.00
44) 2-Fluorobiphenyl	12.95	172	3997220	81.32	ng	0.00
78) Terphenyl-d14	19.71	244	4044445	75.78	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.29	88	261151	39.254	ng	# 78
3) Pyridine	3.66	79	776665	40.302	ng	# 72
4) n-Nitrosodimethylamine	3.59	42	314201	38.009	ng	# 73
6) Aniline	7.03	93	1153338	37.966	ng	98
8) 2-Chlorophenol	7.26	128	776610	37.788	ng	96
9) Benzaldehyde	6.85	77	593741	38.767	ng	91
10) Phenol	6.90	94	974144	38.215	ng	98
11) bis(2-Chloroethyl)ether	7.13	93	789282	38.041	ng	96
12) 1,3-Dichlorobenzene	7.58	146	891872	38.079	ng	99
13) 1,4-Dichlorobenzene	7.73	146	913684	38.132	ng	97
14) 1,2-Dichlorobenzene	8.04	146	869929	37.635	ng	97
15) Benzyl Alcohol	7.93	79	701480	37.381	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.22	45	1275127	38.059	ng	94
17) 2-Methylphenol	8.13	107	644536	37.036	ng	95
18) Hexachloroethane	8.76	117	334081	38.096	ng	90
19) n-Nitroso-di-n-propylamine	8.50	70	653219	36.213	ng	# 97
20) 3+4-Methylphenols	8.46	107	893152	36.784	ng	96
22) Acetophenone	8.51	105	1190230	37.942	ng	96
24) Nitrobenzene	8.89	77	939482	38.424	ng	94
25) Isophorone	9.40	82	1488703	36.750	ng	97
26) 2-Nitrophenol	9.59	139	450544	38.146	ng	97
27) 2,4-Dimethylphenol	9.65	122	619904	37.739	ng	96
28) bis(2-Chloroethoxy)methane	9.89	93	978496	37.395	ng	96
29) 2,4-Dichlorophenol	10.12	162	732450	38.239	ng	94
30) 1,2,4-Trichlorobenzene	10.33	180	824959	38.030	ng	98
31) Naphthalene	10.52	128	2300033	37.668	ng	98
32) Benzoic acid	9.82	122	445250	31.948	ng	93
33) 4-Chloroaniline	10.63	127	1000018	36.808	ng	98
34) Hexachlorobutadiene	10.80	225	509484	38.353	ng	97
35) Caprolactam	11.41	113	215397	32.951	ng	# 66
36) 4-Chloro-3-methylphenol	11.76	107	783793	35.986	ng	95
37) 2-Methylnaphthalene	12.13	142	1571055	36.481	ng	95
39) 1,2,4,5-Tetrachlorobenzene	12.50	216	862257	40.055	ng	98
40) Hexachlorocyclopentadiene	12.48	237	518383	42.637	ng	92

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.75	196	560746	38.824	nq	98
43) 2,4,5-Trichlorophenol	12.82	196	577597	37.752	nq	98
45) 1,1'-Biphenyl	13.15	154	2190035	38.874	nq	98
46) 2-Chloronaphthalene	13.19	162	1702420	39.150	nq	97
47) 2-Nitroaniline	13.40	65	532531	38.059	nq	92
48) Acenaphthylene	14.05	152	2504558	37.955	nq	98
49) Dimethylphthalate	13.79	163	1907982	35.878	nq	99
50) 2,6-Dinitrotoluene	13.90	165	435137	36.830	nq	95
51) Acenaphthene	14.39	154	1494309	37.594	nq	98
52) 3-Nitroaniline	14.23	138	464482	36.284	nq	# 87
53) 2,4-Dinitrophenol	14.45	184	208876	33.389	nq	96
54) Dibenzofuran	14.73	168	2377555	36.893	nq	97
55) 4-Nitrophenol	14.55	139	344846	35.088	nq	91
56) 2,4-Dinitrotoluene	14.70	165	574789	35.846	nq	88
57) Fluorene	15.37	166	1752278	36.075	nq	98
58) 2,3,4,6-Tetrachlorophenol	14.96	232	475815	36.162	nq	94
59) Diethylphthalate	15.16	149	1892419	34.962	nq	99
60) 4-Chlorophenyl-phenylether	15.37	204	983175	36.579	nq	98
61) 4-Nitroaniline	15.40	138	459649	35.224	nq	93
62) Azobenzene	15.67	77	1902404	36.629	nq	99
64) 4,6-Dinitro-2-methylphenol	15.46	198	325741	38.525	nq	98
65) n-Nitrosodiphenylamine	15.59	169	1631199	39.251	nq	97
66) 4-Bromophenyl-phenylether	16.27	248	568256	38.759	nq	# 92
67) Hexachlorobenzene	16.38	284	623917	38.474	nq	98
68) Atrazine	16.55	200	514323	36.759	nq	97
69) Pentachlorophenol	16.73	266	396992	38.031	nq	98
70) Phenanthrene	17.12	178	2590826	37.900	nq	99
71) Anthracene	17.20	178	2553221	37.735	nq	98
72) Carbazole	17.47	167	2514490	37.037	nq	97
73) Di-n-butylphthalate	18.05	149	2920916	36.883	nq	99
74) Fluoranthene	19.13	202	2738932	36.903	nq	98
76) Benzidine	19.32	184	1323448	36.456	nq	96
77) Pyrene	19.50	202	2810577	36.615	nq	100
79) Butylbenzylphthalate	20.41	149	1230154	36.816	nq	93
80) Benzo(a)anthracene	21.25	228	2526342	37.239	nq	98
81) 3,3'-Dichlorobenzidine	21.19	252	1017610	38.502	nq	98
82) Chrysene	21.30	228	2437439	37.735	nq	98
83) Bis(2-ethylhexyl)phthalate	21.19	149	1749068	37.538	nq	96
84) Di-n-octyl phthalate	22.07	149	2901692	39.276	nq	97
85) Indeno(1,2,3-cd)pyrene	25.73	276	2723988	43.254	nq	# 95
87) Benzo(b)fluoranthene	22.83	252	2383242	36.730	nq	98
88) Benzo(k)fluoranthene	22.87	252	2394940m	37.479	nq	
89) Benzo(a)pyrene	23.40	252	2306889	38.070	nq	# 97
90) Dibenzo(a,h)anthracene	25.74	278	2292960	40.304	nq	# 95
91) Benzo(g,h,i)perylene	26.40	276	2234175	40.635	nq	# 95

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

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