

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072518\
 Data File : BN002132.D
 Acq On : 25 Jul 2018 12:53
 Operator : JU/SJ
 Sample : SSTDICC2.5
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC2.5

Manual Integrations
 APPROVED

Sohil
 7/27/2018 11:35:21 AM

Quant Time: Jul 25 20:03:04 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 19:40:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	213689	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	958961	20.00	ng	0.00
38) Acenaphthene-d10	14.32	164	601798	20.00	ng	0.00
63) Phenanthrene-d10	17.07	188	1392975	20.00	ng	0.00
75) Chrysene-d12	21.27	240	1345358	20.00	ng	0.00
86) Perylene-d12	23.49	264	1184178	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.31	112	62146	4.68	ng	0.00
7) Phenol-d6	6.88	99	83561	4.49	ng	0.00
23) Nitrobenzene-d5	8.84	82	86144	4.61	ng	0.00
41) 2,4,6-Tribromophenol	15.82	330	36882	4.72	ng	0.00
44) 2-Fluorobiphenyl	12.94	172	223999	4.85	ng	0.00
78) Terphenyl-d14	19.70	244	310797	4.82	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.29	88	12941	2.597	ng	# 84
3) Pyridine	3.68	79	36214m	2.509	ng	
4) n-Nitrosodimethylamine	3.59	42	17704m	2.859	ng	
6) Aniline	7.03	93	51852	2.279	ng	97
8) 2-Chlorophenol	7.26	128	35738	2.322	ng	92
9) Benzaldehyde	6.85	77	30616	2.657	ng	94
10) Phenol	6.90	94	42828	2.243	ng	99
11) bis(2-Chloroethyl)ether	7.13	93	37908	2.439	ng	93
12) 1,3-Dichlorobenzene	7.58	146	42330	2.413	ng	96
13) 1,4-Dichlorobenzene	7.73	146	43747	2.438	ng	98
14) 1,2-Dichlorobenzene	8.04	146	43654	2.521	ng	96
15) Benzyl Alcohol	7.93	79	35195m	2.504	ng	
16) 2,2'-oxybis(1-Chloropropan	8.22	45	62269	2.481	ng	92
17) 2-Methylphenol	8.13	107	31033	2.381	ng	97
18) Hexachloroethane	8.76	117	15895	2.420	ng	92
19) n-Nitroso-di-n-propylamine	8.49	70	31862	2.358	ng	# 96
20) 3+4-Methylphenols	8.46	107	41231	2.267	ng	94
22) Acetophenone	8.50	105	58806	2.365	ng	93
24) Nitrobenzene	8.88	77	45594	2.353	ng	# 97
25) Isophorone	9.40	82	73922	2.302	ng	96
26) 2-Nitrophenol	9.59	139	20329	2.172	ng	90
27) 2,4-Dimethylphenol	9.65	122	28587	2.196	ng	94
28) bis(2-Chloroethoxy)methane	9.89	93	49696	2.396	ng	# 97
29) 2,4-Dichlorophenol	10.12	162	33431	2.202	ng	94
30) 1,2,4-Trichlorobenzene	10.33	180	41618	2.421	ng	96
31) Naphthalene	10.52	128	119405	2.467	ng	99
33) 4-Chloroaniline	10.63	127	47980	2.228	ng	98
34) Hexachlorobutadiene	10.80	225	24967	2.371	ng	97
35) Caprolactam	11.38	113	10911	2.106	ng	# 63
36) 4-Chloro-3-methylphenol	11.76	107	38217	2.214	ng	# 85
37) 2-Methylnaphthalene	12.13	142	83147	2.436	ng	92
39) 1,2,4,5-Tetrachlorobenzene	12.50	216	47260	2.336	ng	97
40) Hexachlorocyclopentadiene	12.48	237	20438	1.788	ng	89
42) 2,4,6-Trichlorophenol	12.75	196	29206	2.151	ng	95

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43) 2,4,5-Trichlorophenol	12.83	196	31259	2.174	nq	99
45) 1,1'-Biphenyl	13.15	154	125802	2.376	nq	97
46) 2-Chloronaphthalene	13.19	162	95902	2.346	nq	97
47) 2-Nitroaniline	13.40	65	27379	2.082	nq	# 88
48) Acenaphthylene	14.04	152	146457	2.361	nq	95
49) Dimethylphthalate	13.78	163	122583	2.452	nq	100
50) 2,6-Dinitrotoluene	13.90	165	24533	2.209	nq	92
51) Acenaphthene	14.39	154	90314	2.417	nq	97
52) 3-Nitroaniline	14.23	138	25341	2.106	nq	# 93
54) Dibenzofuran	14.72	168	150106	2.478	nq	95
55) 4-Nitrophenol	14.56	139	17948	1.943	nq	90
56) 2,4-Dinitrotoluene	14.70	165	33237	2.205	nq	88
57) Fluorene	15.37	166	113041	2.476	nq	98
58) 2,3,4,6-Tetrachlorophenol	14.96	232	28934	2.339	nq	# 95
59) Diethylphthalate	15.15	149	125450	2.466	nq	98
60) 4-Chlorophenyl-phenylether	15.37	204	63321	2.506	nq	98
61) 4-Nitroaniline	15.40	138	26075	2.126	nq	90
62) Azobenzene	15.66	77	117933	2.416	nq	98
64) 4,6-Dinitro-2-methylphenol	15.46	198	16265	1.729	nq	99
65) n-Nitrosodiphenylamine	15.59	169	105577	2.284	nq	96
66) 4-Bromophenyl-phenylether	16.27	248	37907	2.325	nq	93
67) Hexachlorobenzene	16.38	284	43277	2.399	nq	95
68) Atrazine	16.54	200	36808	2.365	nq	96
69) Pentachlorophenol	16.73	266	20608	1.775	nq	97
70) Phenanthrene	17.11	178	186991	2.459	nq	98
71) Anthracene	17.20	178	178941	2.378	nq	97
72) Carbazole	17.47	167	180279	2.387	nq	96
73) Di-n-butylphthalate	18.05	149	200736	2.279	nq	98
74) Fluoranthene	19.13	202	206442	2.501	nq	97
76) Benzidine	19.32	184	104636	2.385	nq	97
77) Pyrene	19.50	202	216327	2.332	nq	99
79) Butylbenzylphthalate	20.41	149	85379	2.114	nq	92
80) Benzo(a)anthracene	21.25	228	199328	2.431	nq	97
81) 3,3'-Dichlorobenzidine	21.19	252	73129	2.289	nq	# 97
82) Chrysene	21.30	228	192527	2.466	nq	96
83) Bis(2-ethylhexyl)phthalate	21.19	149	120267	2.135	nq	91
84) Di-n-octyl phthalate	22.07	149	183328	2.053	nq	100
85) Indeno(1,2,3-cd)pyrene	25.72	276	160421	2.583	nq	# 94
87) Benzo(b)fluoranthene	22.83	252	172540	2.469	nq	98
88) Benzo(k)fluoranthene	22.87	252	166340	2.417	nq	98
89) Benzo(a)pyrene	23.40	252	156175	2.393	nq	97
90) Dibenzo(a,h)anthracene	25.73	278	136045	2.220	nq	# 93
91) Benzo(g,h,i)perylene	26.40	276	128301	2.167	nq	# 96

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

