

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050818\
 Data File : BF105140.D
 Acq On : 8 May 2018 12:06
 Operator : JU/SJ
 Sample : SSTDICC025
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC025

Manual Integrations
 APPROVED

Sohil
 5/9/2018 6:50:34 PM

Quant Time: May 08 14:49:24 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 08 13:40:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	200472	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	802335	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	396320	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	756711	20.00	ng	0.00
75) Chrysene-d12	14.03	240	621221	20.00	ng	0.01
86) Perylene-d12	15.60	264	590536	20.00	ng	0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.39	112	639727	48.79	ng	0.00
7) Phenol-d6	6.41	99	748452	46.46	ng	-0.01
23) Nitrobenzene-d5	7.36	82	591862	48.17	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	223744	54.42	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1349562	53.79	ng	0.00
78) Terphenyl-d14	12.91	244	1360301	49.92	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.52	88	157320	25.752	ng	87
3) Pyridine	3.24	79	438261	25.516	ng	86
4) n-Nitrosodimethylamine	3.19	42	219609	36.576	ng	84
6) Aniline	6.46	93	537172	24.002	ng	97
8) 2-Chlorophenol	6.57	128	339930	24.565	ng	97
9) Benzaldehyde	6.35	77	295190	36.722	ng	99
10) Phenol	6.43	94	417013	24.398	ng	97
11) bis(2-Chloroethyl)ether	6.53	93	337930	24.776	ng	91
12) 1,3-Dichlorobenzene	6.73	146	393577	25.105	ng	100
13) 1,4-Dichlorobenzene	6.81	146	396910	25.398	ng	99
14) 1,2-Dichlorobenzene	6.96	146	374526	25.862	ng	98
15) Benzyl Alcohol	6.93	79	287500	23.498	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.07	45	632120	34.509	ng	96
17) 2-Methylphenol	7.04	107	287292	23.884	ng	99
18) Hexachloroethane	7.30	117	126773	22.753	ng	97
19) n-Nitroso-di-n-propylamine	7.20	70	248249	23.583	ng	90
20) 3+4-Methylphenols	7.19	107	354408	23.756	ng	89
22) Acetophenone	7.20	105	482969	24.450	ng	# 96
24) Nitrobenzene	7.37	77	342515	25.248	ng	95
25) Isophorone	7.62	82	676754	26.046	ng	99
26) 2-Nitrophenol	7.69	139	86682	15.695	ng	99
27) 2,4-Dimethylphenol	7.72	122	311599	27.173	ng	98
28) bis(2-Chloroethoxy)methane	7.83	93	423181	26.638	ng	98
29) 2,4-Dichlorophenol	7.93	162	284004	25.957	ng	94
30) 1,2,4-Trichlorobenzene	8.02	180	325125	27.076	ng	98
31) Naphthalene	8.10	128	1000553	25.044	ng	100
32) Benzoic acid	7.82	122	96815m	10.581	ng	
33) 4-Chloroaniline	8.15	127	423489	24.742	ng	98
34) Hexachlorobutadiene	8.22	225	190068	28.898	ng	99
35) Caprolactam	8.50	113	83053	21.759	ng	# 79
36) 4-Chloro-3-methylphenol	8.62	107	296273	25.253	ng	99
37) 2-Methylnaphthalene	8.79	142	707696	27.385	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.95	216	331017	29.177	ng	98
40) Hexachlorocyclopentadiene	8.94	237	162705	25.618	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.06	196	196832	24.993	nq	95
43) 2,4,5-Trichlorophenol	9.10	196	207330	23.526	nq #	90
45) 1,1'-Biphenyl	9.25	154	853188	25.626	nq	98
46) 2-Chloronaphthalene	9.27	162	628142	23.886	nq	97
47) 2-Nitroaniline	9.37	65	146945	22.595	nq	85
48) Acenaphthylene	9.69	152	981068	23.778	nq	99
49) Dimethylphthalate	9.55	163	703387	22.506	nq	98
50) 2,6-Dinitrotoluene	9.62	165	143311	24.781	nq	93
51) Acenaphthene	9.86	154	610778	24.262	nq	97
52) 3-Nitroaniline	9.78	138	151970	22.447	nq	99
53) 2,4-Dinitrophenol	9.88	184	23220	13.209	nq #	20
54) Dibenzofuran	10.03	168	921839	26.276	nq	100
55) 4-Nitrophenol	9.93	139	115914	21.796	nq	94
56) 2,4-Dinitrotoluene	10.01	165	173681	22.896	nq	92
57) Fluorene	10.37	166	675074	26.436	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.14	232	185187	28.166	nq #	88
59) Diethylphthalate	10.25	149	704722	22.641	nq	96
60) 4-Chlorophenyl-phenylether	10.37	204	325811	28.649	nq	89
61) 4-Nitroaniline	10.39	138	148308	22.404	nq	94
62) Azobenzene	10.53	77	710130	23.880	nq	92
64) 4,6-Dinitro-2-methylphenol	10.42	198	42408	12.900	nq	93
65) n-Nitrosodiphenylamine	10.49	169	612245	23.335	nq	100
66) 4-Bromophenyl-phenylether	10.86	248	212873	26.447	nq #	85
67) Hexachlorobenzene	10.92	284	242230	26.746	nq #	87
68) Atrazine	11.02	200	196433	24.803	nq	98
69) Pentachlorophenol	11.12	266	123542	22.049	nq	98
70) Phenanthrene	11.34	178	1045961	24.821	nq	99
71) Anthracene	11.39	178	1085358	25.337	nq	99
72) Carbazole	11.54	167	970922	23.195	nq	98
73) Di-n-butylphthalate	11.88	149	1053956	20.612	nq	99
74) Fluoranthene	12.53	202	1117195	25.374	nq	98
76) Benzidine	12.66	184	614090	28.559	nq	99
77) Pyrene	12.76	202	1137169	24.696	nq	100
79) Butylbenzylphthalate	13.41	149	316675	14.598	nq	98
80) Benzo(a)anthracene	14.02	228	980906	26.431	nq	99
81) 3,3'-Dichlorobenzidine	13.98	252	363218	26.249	nq #	96
82) Chrysene	14.06	228	991277	26.004	nq	99
83) Bis(2-ethylhexyl)phthalate	14.02	149	474111	16.991	nq	100
84) Di-n-octyl phthalate	14.72	149	724148	15.532	nq #	92
85) Indeno(1,2,3-cd)pyrene	17.04	276	770075	23.781	nq	94
87) Benzo(b)fluoranthene	15.17	252	955624	25.622	nq	98
88) Benzo(k)fluoranthene	15.20	252	903195	25.650	nq	98
89) Benzo(a)pyrene	15.53	252	870720	26.092	nq	98
90) Dibenzo(a,h)anthracene	17.07	278	638447	23.294	nq	96
91) Benzo(g,h,i)perylene	17.48	276	626219	23.583	nq	95

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

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