

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072717\
 Data File : BF097118.D
 Acq On : 27 Jul 2017 13:17
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jul 27 15:52:35 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 14:11:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.49	152	112994	20.00	ng	-0.01
21) Naphthalene-d8	7.77	136	478336	20.00	ng	-0.01
38) Acenaphthene-d10	9.52	164	210077	20.00	ng	-0.02
63) Phenanthrene-d10	10.99	188	370199	20.00	ng	-0.01
75) Chrysene-d12	13.62	240	227072	20.00	ng	-0.01
86) Perylene-d12	14.96	264	206145	20.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.08	112	583844	77.89	ng	-0.02
7) Phenol-d6	6.15	99	674091	78.78	ng	-0.02
23) Nitrobenzene-d5	7.06	82	606153	74.34	ng	-0.01
41) 2,4,6-Tribromophenol	10.30	330	138644	83.53	ng	-0.02
44) 2-Fluorobiphenyl	8.85	172	959941	77.55	ng	-0.02
78) Terphenyl-d14	12.57	244	875856	78.64	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.01	88	126910	40.573	ng	# 77
3) Pyridine	2.63	79	417835	43.624	ng	# 66
4) n-Nitrosodimethylamine	2.59	42	229659	43.281	ng	# 80
6) Aniline	6.15	93	415102	38.549	ng	# 79
8) 2-Chlorophenol	6.27	128	324512	40.489	ng	# 92
9) Benzaldehyde	6.03	77	190528	37.617	ng	# 99
10) Phenol	6.17	94	403078	39.712	ng	# 94
11) bis(2-Chloroethyl)ether	6.23	93	303612	37.821	ng	# 91
12) 1,3-Dichlorobenzene	6.43	146	331419	38.371	ng	# 98
13) 1,4-Dichlorobenzene	6.50	146	337179	39.391	ng	# 95
14) 1,2-Dichlorobenzene	6.66	146	302675	40.498	ng	# 99
15) Benzyl Alcohol	6.65	79	213580	39.423	ng	# 96
16) 2,2'-oxybis(1-Chloropropan	6.78	45	596050	37.019	ng	# 93
17) 2-Methylphenol	6.77	107	237842	38.450	ng	# 90
18) Hexachloroethane	7.00	117	123899	39.565	ng	# 83
19) n-Nitroso-di-n-propylamine	6.92	70	208241	36.971	ng	# 83
20) 3+4-Methylphenols	6.93	107	316135	39.130	ng	# 63
22) Acetophenone	6.91	105	426418	37.122	ng	# 97
24) Nitrobenzene	7.08	77	321030	37.804	ng	# 91
25) Isophorone	7.32	82	590879	37.003	ng	# 97
26) 2-Nitrophenol	7.39	139	175403	38.178	ng	# 85
27) 2,4-Dimethylphenol	7.45	122	286511	37.804	ng	# 97
28) bis(2-Chloroethoxy)methane	7.54	93	369882	35.495	ng	# 99
29) 2,4-Dichlorophenol	7.65	162	252795	38.719	ng	# 97
30) 1,2,4-Trichlorobenzene	7.72	180	246016	39.532	ng	# 97
31) Naphthalene	7.79	128	867101	38.455	ng	# 99
32) Benzoic acid	7.62	122	226408	40.007	ng	# 94
33) 4-Chloroaniline	7.85	127	357860	39.005	ng	# 99
34) Hexachlorobutadiene	7.92	225	127052	38.510	ng	# 98
35) Caprolactam	8.24	113	91459	43.686	ng	# 53
36) 4-Chloro-3-methylphenol	8.35	107	276123	38.765	ng	# 87
37) 2-Methylnaphthalene	8.49	142	547606	38.357	ng	# 98
39) 1,2,4,5-Tetrachlorobenzene	8.66	216	227076	40.510	ng	# 99
40) Hexachlorocyclopentadiene	8.64	237	53960	31.852	ng	# 98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.77	196	157416	38.753	nq	98
43) 2,4,5-Trichlorophenol	8.82	196	157839	38.821	nq	98
45) 1,1'-Biphenyl	8.95	154	693170	38.631	nq	99
46) 2-Chloronaphthalene	8.97	162	504381	38.198	nq	# 92
47) 2-Nitroaniline	9.07	65	179876	37.536	nq	95
48) Acenaphthylene	9.38	152	762316	37.612	nq	99
49) Dimethylphthalate	9.26	163	632858	39.670	nq	99
50) 2,6-Dinitrotoluene	9.32	165	131880	38.977	nq	# 87
51) Acenaphthene	9.55	154	451633	37.830	nq	98
52) 3-Nitroaniline	9.49	138	167136	39.797	nq	# 88
53) 2,4-Dinitrophenol	9.60	184	65178	42.367	nq	# 75
54) Dibenzofuran	9.72	168	613121	38.557	nq	96
55) 4-Nitrophenol	9.67	139	100262	40.144	nq	# 64
56) 2,4-Dinitrotoluene	9.72	165	154156	39.633	nq	# 47
57) Fluorene	10.06	166	489349	40.944	nq	98
58) 2,3,4,6-Tetrachlorophenol	9.85	232	111832	39.836	nq	98
59) Diethylphthalate	9.96	149	568432	38.839	nq	99
60) 4-Chlorophenyl-phenylether	10.06	204	202551	41.041	nq	99
61) 4-Nitroaniline	10.10	138	166738	41.783	nq	93
62) Azobenzene	10.22	77	524634	38.640	nq	92
64) 4,6-Dinitro-2-methylphenol	10.13	198	93302	40.559	nq	91
65) n-Nitrosodiphenylamine	10.18	169	468144	36.345	nq	97
66) 4-Bromophenyl-phenylether	10.55	248	132078	35.750	nq	96
67) Hexachlorobenzene	10.62	284	150992	36.445	nq	96
68) Atrazine	10.72	200	139740	37.833	nq	94
69) Pentachlorophenol	10.81	266	73126	42.659	nq	98
70) Phenanthrene	11.02	178	744680	37.543	nq	99
71) Anthracene	11.07	178	762441	37.530	nq	100
72) Carbazole	11.23	167	743268	37.200	nq	99
73) Di-n-butylphthalate	11.57	149	899179	36.407	nq	99
74) Fluoranthene	12.20	202	715001	36.795	nq	97
76) Benzidine	12.32	184	355217	42.160	nq	# 94
77) Pyrene	12.42	202	728953	39.213	nq	99
79) Butylbenzylphthalate	13.05	149	363016	38.390	nq	# 77
80) Benzo(a)anthracene	13.60	228	567503	38.625	nq	98
81) 3,3'-Dichlorobenzidine	13.57	252	226373	40.857	nq	# 96
82) Chrysene	13.64	228	554734	38.666	nq	99
83) Bis(2-ethylhexyl)phthalate	13.62	149	449223	36.385	nq	100
84) Di-n-octyl phthalate	14.23	149	774135	34.167	nq	# 93
85) Indeno(1,2,3-cd)pyrene	16.19	276	429552	34.917	nq	99
87) Benzo(b)fluoranthene	14.60	252	508978	41.074	nq	98
88) Benzo(k)fluoranthene	14.63	252	453322	38.390	nq	97
89) Benzo(a)pyrene	14.91	252	429167	37.841	nq	99
90) Dibenzo(a,h)anthracene	16.19	278	345564	37.156	nq	96
91) Benzo(g,h,i)perylene	16.54	276	378577	38.816	nq	99

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

