

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050916\
 Data File : BF086845.D
 Acq On : 10 May 2016 2:51
 Operator : UM/SJ
 Sample : H2895-11MSD
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP8-9.5FTMSD

Manual Integrations
 APPROVED

sohil
 5/10/2016 11:06:08 AM

Quant Time: May 10 03:58:45 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 09 16:04:56 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	171989	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	719308	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	337907	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	595981	20.00	ng	0.00
75) Chrysene-d12	13.84	240	340753	20.00	ng	0.00
86) Perylene-d12	15.21	264	347559	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	1262725	124.34	ng	0.01
7) Phenol-d6	6.36	99	1654303	121.88	ng	0.00
23) Nitrobenzene-d5	7.26	82	1194544	83.39	ng	0.00
41) 2,4,6-Tribromophenol	10.52	330	411462	115.02	ng	0.00
44) 2-Fluorobiphenyl	9.06	172	1565487	77.86	ng	0.00
78) Terphenyl-d14	12.79	244	1328832	82.74	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.25	88	152696	30.63	ng	# 69
3) Pyridine	2.92	79	432381	35.47	ng	# 67
4) n-Nitrosodimethylamine	2.88	42	348284	39.40	ng	# 51
6) Aniline	6.36	93	437530	26.65	ng	# 45
8) 2-Chlorophenol	6.49	128	499734	40.78	ng	97
9) Benzaldehyde	6.24	77	49471	6.29	ng	92
10) Phenol	6.37	94	711129	44.08	ng	72
11) bis(2-Chloroethyl)ether	6.44	93	502985	39.99	ng	91
12) 1,3-Dichlorobenzene	6.62	146	491786m	37.08	ng	
13) 1,4-Dichlorobenzene	6.70	146	494295m	36.55	ng	
14) 1,2-Dichlorobenzene	6.86	146	476679	40.44	ng	99
15) Benzyl Alcohol	6.85	79	430566	45.94	ng	# 87
16) 2,2'-oxybis(1-Chloropropan	6.98	45	982760	35.93	ng	56
17) 2-Methylphenol	6.98	107	424509	43.53	ng	# 82
18) Hexachloroethane	7.20	117	186262	36.61	ng	95
19) n-Nitroso-di-n-propylamine	7.13	70	429316	45.26	ng	# 78
20) 3+4-Methylphenols	7.14	107	566381	44.47	ng	# 58
22) Acetophenone	7.12	105	730259	42.98	ng	99
24) Nitrobenzene	7.29	77	618548	41.97	ng	98
25) Isophorone	7.53	82	1049239	39.48	ng	99
26) 2-Nitrophenol	7.61	139	268248	41.41	ng	94
27) 2,4-Dimethylphenol	7.65	122	433346	39.27	ng	86
28) bis(2-Chloroethoxy)methane	7.74	93	669813	46.72	ng	98
29) 2,4-Dichlorophenol	7.85	162	405553	41.76	ng	92
30) 1,2,4-Trichlorobenzene	7.93	180	447864	41.24	ng	98
31) Naphthalene	8.00	128	1429163	39.67	ng	99
32) Benzoic acid	7.77	122	66416	10.25	ng	# 82
33) 4-Chloroaniline	8.06	127	248553	17.75	ng	96
34) Hexachlorobutadiene	8.12	225	248985	39.52	ng	98
35) Caprolactam	8.45	113	140812m	42.61	ng	
36) 4-Chloro-3-methylphenol	8.57	107	458506	38.93	ng	91
37) 2-Methylnaphthalene	8.69	142	934090	44.35	ng	96
39) 1,2,4,5-Tetrachlorobenzene	8.86	216	385084	39.20	ng	99
40) Hexachlorocyclopentadiene	8.84	237	212156	45.50	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.98	196	270334	41.15	nq	96
43) 2,4,5-Trichlorophenol	9.04	196	282254	41.54	nq	95
45) 1,1'-Biphenyl	9.16	154	1186001	44.13	nq	99
46) 2-Chloronaphthalene	9.18	162	886625	42.11	nq	99
47) 2-Nitroaniline	9.29	65	321259	41.74	nq	# 76
48) Acenaphthylene	9.60	152	1321499	42.67	nq	99
49) Dimethylphthalate	9.47	163	1388388	53.86	nq	99
50) 2,6-Dinitrotoluene	9.53	165	223489	41.19	nq	# 65
51) Acenaphthene	9.77	154	906941	47.07	nq	98
52) 3-Nitroaniline	9.70	138	133359	23.35	nq	# 79
53) 2,4-Dinitrophenol	9.81	184	88419	36.52	nq	90
54) Dibenzofuran	9.94	168	1215572	49.15	nq	96
55) 4-Nitrophenol	9.89	139	325175	85.07	nq	# 64
56) 2,4-Dinitrotoluene	9.94	165	292357	43.54	nq	# 87
57) Fluorene	10.28	166	1008432	47.81	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.06	232	241755	47.26	nq	# 100
59) Diethylphthalate	10.17	149	1039789	41.39	nq	97
60) 4-Chlorophenyl-phenylether	10.27	204	372614	39.34	nq	98
61) 4-Nitroaniline	10.32	138	203676	37.13	nq	# 72
62) Azobenzene	10.43	77	1164393	42.96	nq	85
64) 4,6-Dinitro-2-methylphenol	10.34	198	101642	27.45	nq	# 24
65) n-Nitrosodiphenylamine	10.40	169	788116	43.04	nq	99
66) 4-Bromophenyl-phenylether	10.76	248	260277	43.07	nq	# 82
67) Hexachlorobenzene	10.82	284	261570	37.04	nq	# 74
68) Atrazine	10.93	200	272376	44.53	nq	95
69) Pentachlorophenol	11.02	266	279442	85.62	nq	98
70) Phenanthrene	11.24	178	2260384	71.07	nq	99
71) Anthracene	11.29	178	1550302	47.66	nq	99
72) Carbazole	11.45	167	1238926	44.31	nq	98
73) Di-n-butylphthalate	11.78	149	1331608	39.00	nq	# 98
74) Fluoranthene	12.42	202	1888856	57.74	nq	97
76) Benzidine	12.55	184	500985	57.67	nq	97
77) Pyrene	12.65	202	1889105	72.41	nq	99
79) Butylbenzylphthalate	13.27	149	471070	42.69	nq	# 72
80) Benzo(a)anthracene	13.83	228	1055056	55.15	nq	100
81) 3,3'-Dichlorobenzidine	13.79	252	163689	13.47	nq	99
82) Chrysene	13.86	228	900863	46.30	nq	98
83) Bis(2-ethylhexyl)phthalate	13.83	149	586001	44.15	nq	97
84) Di-n-octyl phthalate	14.43	149	979227	44.54	nq	# 85
85) Indeno(1,2,3-cd)pyrene	16.51	276	962716	59.46	nq	95
87) Benzo(b)fluoranthene	14.83	252	1035184m	45.83	nq	
88) Benzo(k)fluoranthene	14.85	252	790875m	42.76	nq	
89) Benzo(a)pyrene	15.16	252	976879	50.14	nq	# 96
90) Dibenzo(a,h)anthracene	16.52	278	763026	46.47	nq	# 92
91) Benzo(g,h,i)perylene	16.90	276	828955	49.68	nq	# 92

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

