

Data Path : Z:\HPCHEM1\BNA F\DATA\BF021416\
 Data File : BF085065.D
 Acq On : 13 Feb 2016 3:48
 Operator : SJ/IZ
 Sample : H1396-07MSD
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S4-B003(10-12)MSD

Manual Integrations
 APPROVED

Sohil
 2/15/2016 4:44:46 PM

Quant Time: Feb 13 06:43:56 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 13 00:11:18 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.88	152	145953	20.00	ng	0.00
21) Naphthalene-d8	7.78	136	654048	20.00	ng	-0.01
38) Acenaphthene-d10	10.58	164	297744	20.00	ng	0.00
63) Phenanthrene-d10	12.99	188	531089	20.00	ng	0.00
75) Chrysene-d12	17.18	240	339258	20.00	ng	0.00
86) Perylene-d12	18.80	264	279888	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	4.15	112	915928	109.73	ng	-0.01
7) Phenol-d6	5.44	99	1386307	124.62	ng	0.00
23) Nitrobenzene-d5	6.71	82	886050	80.00	ng	-0.01
41) 2,4,6-Tribromophenol	11.88	330	347520	113.00	ng	0.00
44) 2-Fluorobiphenyl	9.53	172	1261843	60.58	ng	0.00
78) Terphenyl-d14	15.62	244	984590	59.73	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.02	88	116337	33.83	ng	99
3) Pyridine	2.47	79	246410	30.67	ng	96
4) n-Nitrosodimethylamine	2.46	42	119282	28.86	ng	# 98
6) Aniline	5.46	93	220349	18.11	ng	# 80
8) 2-Chlorophenol	5.59	128	418853	40.82	ng	95
9) Benzaldehyde	5.42	77	107126	17.88	ng	92
10) Phenol	5.46	94	503579	38.70	ng	85
11) bis(2-Chloroethyl)ether	5.53	93	436712	38.57	ng	98
12) 1,3-Dichlorobenzene	5.79	146	378041	37.05	ng	95
13) 1,4-Dichlorobenzene	5.91	146	388876	34.25	ng	97
14) 1,2-Dichlorobenzene	6.11	146	425329	39.85	ng	99
15) Benzyl Alcohol	6.19	79	220424	32.43	ng	97
16) 2,2'-oxybis(1-Chloropropan	6.30	45	677048	39.50	ng	94
17) 2-Methylphenol	6.33	107	243378	32.76	ng	95
18) Hexachloroethane	6.58	117	168735	40.33	ng	90
19) n-Nitroso-di-n-propylamine	6.50	70	286222	40.84	ng	95
20) 3+4-Methylphenols	6.57	107	290554	32.48	ng	# 81
22) Acetophenone	6.50	105	607204	36.22	ng	# 96
24) Nitrobenzene	6.74	77	335648	27.74	ng	99
25) Isophorone	7.10	82	785978	39.56	ng	97
26) 2-Nitrophenol	7.23	139	223030	38.45	ng	98
27) 2,4-Dimethylphenol	7.35	122	366341	39.03	ng	97
28) bis(2-Chloroethoxy)methane	7.47	93	496439	40.25	ng	99
29) 2,4-Dichlorophenol	7.63	162	322718	37.83	ng	99
30) 1,2,4-Trichlorobenzene	7.70	180	381052	38.75	ng	99
31) Naphthalene	7.82	128	1271719	40.10	ng	99
33) 4-Chloroaniline	8.03	127	127779	10.62	ng	96
34) Hexachlorobutadiene	8.01	225	208607	37.88	ng	96
35) Caprolactam	8.57	113	80932m	35.29	ng	
36) 4-Chloro-3-methylphenol	8.81	107	359001	37.12	ng	97
37) 2-Methylnaphthalene	8.92	142	792175	39.16	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.19	216	365128	37.95	ng	98
40) Hexachlorocyclopentadiene	9.16	237	249152	64.35	ng	97
42) 2,4,6-Trichlorophenol	9.43	196	187631	35.37	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	9.51	196	249540	40.43	nq	96
45) 1,1'-Biphenyl	9.69	154	1036019	38.25	nq	99
46) 2-Chloronaphthalene	9.70	162	762138	39.71	nq	100
47) 2-Nitroaniline	9.94	65	232961	39.32	nq	97
48) Acenaphthylene	10.35	152	1159353	38.05	nq	99
49) Dimethylphthalate	10.23	163	1112098	49.82	nq	99
50) 2,6-Dinitrotoluene	10.33	165	175716	38.37	nq	# 90
51) Acenaphthene	10.64	154	710804	38.58	nq	99
52) 3-Nitroaniline	10.66	138	123687	24.31	nq	99
53) 2,4-Dinitrophenol	10.81	184	81859	45.91	nq	98
54) Dibenzofuran	10.92	168	1039094	42.17	nq	99
55) 4-Nitrophenol	11.04	139	209604	64.20	nq	91
56) 2,4-Dinitrotoluene	10.99	165	251947	40.38	nq	93
57) Fluorene	11.48	166	828285	39.19	nq	99
58) 2,3,4,6-Tetrachlorophenol	11.16	232	227519	46.60	nq	# 97
59) Diethylphthalate	11.38	149	865981	39.15	nq	98
60) 4-Chlorophenyl-phenylether	11.51	204	392879	38.80	nq	95
61) 4-Nitroaniline	11.64	138	168362	36.01	nq	97
62) Azobenzene	11.76	77	898532	42.16	nq	97
64) 4,6-Dinitro-2-methylphenol	11.64	198	100228	34.33	nq	85
65) n-Nitrosodiphenylamine	11.72	169	684835	40.61	nq	99
66) 4-Bromophenyl-phenylether	12.30	248	240539	40.66	nq	95
67) Hexachlorobenzene	12.35	284	256855	40.61	nq	92
68) Atrazine	12.64	200	180309	36.89	nq	97
69) Pentachlorophenol	12.72	266	194811	71.67	nq	97
70) Phenanthrene	13.03	178	1118667	38.20	nq	98
71) Anthracene	13.12	178	1268186	41.44	nq	99
72) Carbazole	13.43	167	1048962	41.40	nq	99
73) Di-n-butylphthalate	14.03	149	1292140	38.85	nq	99
74) Fluoranthene	14.96	202	1121511	38.44	nq	100
76) Benzidine	15.28	184	155776	22.55	nq	98
77) Pyrene	15.31	202	1111714	40.97	nq	98
79) Butylbenzylphthalate	16.45	149	460118	40.32	nq	99
80) Benzo(a)anthracene	17.17	228	755223	38.44	nq	99
81) 3,3'-Dichlorobenzidine	17.18	252	103941	18.21	nq	99
82) Chrysene	17.21	228	735721	38.62	nq	98
83) Bis(2-ethylhexyl)phthalate	17.27	149	599572	39.96	nq	99
84) Di-n-octyl phthalate	18.02	149	882769	41.41	nq	99
85) Indeno(1,2,3-cd)pyrene	20.09	276	708357	44.75	nq	99
87) Benzo(b)fluoranthene	18.41	252	642830m	35.03	nq	
88) Benzo(k)fluoranthene	18.43	252	604204m	41.22	nq	
89) Benzo(a)pyrene	18.74	252	595756	38.39	nq	99
90) Dibenzo(a,h)anthracene	20.11	278	603590	40.98	nq	97
91) Benzo(g,h,i)perylene	20.48	276	604880	39.55	nq	97

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

