

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040517\
 Data File : BF094192.D
 Acq On : 5 Apr 2017 16:34
 Operator : SJ/MA
 Sample : I2523-15MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CEMETERY-3(0-7)MS

Manual Integrations
 APPROVED

mohammad
 4/6/2017 3:38:31 PM

Quant Time: Apr 06 06:41:14 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 31 14:38:29 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.86	152	169748	20.00	ng	-0.04
21) Naphthalene-d8	7.37	136	687290	20.00	ng	-0.04
38) Acenaphthene-d10	9.51	164	316428	20.00	ng	-0.04
63) Phenanthrene-d10	11.33	188	565838	20.00	ng	-0.05
75) Chrysene-d12	14.58	240	443851	20.00	ng	-0.05
86) Perylene-d12	16.20	264	313825	20.00	ng	-0.05

System Monitoring Compounds						
5) 2-Fluorophenol	4.40	112	1060086	105.48	ng	-0.03
7) Phenol-d6	5.47	99	1336409	107.49	ng	-0.04
23) Nitrobenzene-d5	6.52	82	832994	69.17	ng	-0.04
41) 2,4,6-Tribromophenol	10.49	330	285447	114.16	ng	-0.04
44) 2-Fluorobiphenyl	8.69	172	1526576	73.58	ng	-0.04
78) Terphenyl-d14	13.31	244	1271388	64.21	ng	-0.04

Target Compounds					Qvalue	
2) 1,4-Dioxane	2.25	88	144319	29.89	ng	99
3) Pyridine	2.74	79	298436	22.68	ng	97
4) n-Nitrosodimethylamine	2.69	42	193557	35.75	ng	96
6) Aniline	5.49	93	473265	27.36	ng	# 34
8) 2-Chlorophenol	5.63	128	425791	38.54	ng	98
9) Benzaldehyde	5.36	77	222699	26.53	ng	96
10) Phenol	5.49	94	518490	36.65	ng	90
11) bis(2-Chloroethyl)ether	5.57	93	398216	36.02	ng	95
12) 1,3-Dichlorobenzene	5.80	146	448820	36.22	ng	100
13) 1,4-Dichlorobenzene	5.89	146	452147	36.03	ng	97
14) 1,2-Dichlorobenzene	6.06	146	421344	36.46	ng	98
15) Benzyl Alcohol	6.03	79	349009	38.35	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.20	45	550854	32.33	ng	96
17) 2-Methylphenol	6.17	107	345820	36.55	ng	98
18) Hexachloroethane	6.45	117	157078	35.61	ng	# 84
19) n-Nitroso-di-n-propylamine	6.35	70	288411	36.09	ng	97
20) 3+4-Methylphenols	6.36	107	453722	39.60	ng	# 64
22) Acetophenone	6.34	105	588067	38.39	ng	# 87
24) Nitrobenzene	6.54	77	431001	36.29	ng	90
25) Isophorone	6.83	82	771088	36.44	ng	# 94
26) 2-Nitrophenol	6.91	139	234234	41.35	ng	96
27) 2,4-Dimethylphenol	6.98	122	374605	38.38	ng	99
28) bis(2-Chloroethoxy)methane	7.09	93	500867	35.89	ng	99
29) 2,4-Dichlorophenol	7.21	162	364480	39.94	ng	99
30) 1,2,4-Trichlorobenzene	7.30	180	353910	37.65	ng	98
31) Naphthalene	7.39	128	1194224	36.14	ng	100
33) 4-Chloroaniline	7.46	127	433269	32.35	ng	# 94
34) Hexachlorobutadiene	7.55	225	175734	36.46	ng	99
35) Caprolactam	7.90	113	110123m	37.84	ng	
36) 4-Chloro-3-methylphenol	8.07	107	378656	39.71	ng	88
37) 2-Methylnaphthalene	8.24	142	835920	37.95	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.45	216	304041	35.12	ng	99
40) Hexachlorocyclopentadiene	8.43	237	287652	71.01	ng	99
42) 2,4,6-Trichlorophenol	8.59	196	238430	38.93	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	8.64	196	247477	37.34	nq	95
45) 1,1'-Biphenyl	8.81	154	935141	36.64	nq	98
46) 2-Chloronaphthalene	8.83	162	742698	37.17	nq	93
47) 2-Nitroaniline	8.96	65	227037	38.31	nq	87
48) Acenaphthylene	9.33	152	1177351	35.46	nq	99
49) Dimethylphthalate	9.20	163	984882	43.79	nq	98
50) 2,6-Dinitrotoluene	9.26	165	199959	38.78	nq	# 86
51) Acenaphthene	9.55	154	700036	36.13	nq	100
52) 3-Nitroaniline	9.47	138	203158	33.55	nq	85
53) 2,4-Dinitrophenol	9.60	184	37320	17.40	nq	# 71
54) Dibenzofuran	9.76	168	1003748	38.06	nq	99
55) 4-Nitrophenol	9.71	139	321469	67.80	nq	# 70
56) 2,4-Dinitrotoluene	9.76	165	242199	37.39	nq	# 69
57) Fluorene	10.19	166	781488	35.73	nq	99
58) 2,3,4,6-Tetrachlorophenol	9.92	232	173915	38.25	nq	98
59) Diethylphthalate	10.07	149	844622	37.99	nq	99
60) 4-Chlorophenyl-phenylether	10.19	204	337446	36.22	nq	89
61) 4-Nitroaniline	10.22	138	225698	38.85	nq	96
62) Azobenzene	10.39	77	818165	38.90	nq	96
64) 4,6-Dinitro-2-methylphenol	10.26	198	57388	16.56	nq	# 68
65) n-Nitrosodiphenylamine	10.34	169	722134	36.90	nq	99
66) 4-Bromophenyl-phenylether	10.79	248	205528	36.93	nq	96
67) Hexachlorobenzene	10.86	284	209892	37.26	nq	# 89
68) Atrazine	11.01	200	213251	36.38	nq	95
69) Pentachlorophenol	11.11	266	134219	38.28	nq	98
70) Phenanthrene	11.36	178	1137053	36.12	nq	98
71) Anthracene	11.43	178	1184925	36.56	nq	98
72) Carbazole	11.63	167	1108528	33.36	nq	98
73) Di-n-butylphthalate	12.08	149	1307659	37.84	nq	99
74) Fluoranthene	12.82	202	1177915	36.55	nq	99
76) Benzidine	12.99	184	353250	20.11	nq	# 92
77) Pyrene	13.10	202	1182443	36.71	nq	99
79) Butylbenzylphthalate	13.92	149	553094	40.30	nq	# 85
80) Benzo(a)anthracene	14.57	228	957579	37.00	nq	98
81) 3,3'-Dichlorobenzidine	14.54	252	263054	28.43	nq	# 96
82) Chrysene	14.62	228	849468	35.37	nq	99
83) Bis(2-ethylhexyl)phthalate	14.63	149	745173	39.80	nq	# 99
84) Di-n-octyl phthalate	15.39	149	1151279	36.80	nq	98
85) Indeno(1,2,3-cd)pyrene	17.53	276	654262	31.45	nq	99
87) Benzo(b)fluoranthene	15.80	252	744507	38.70	nq	100
88) Benzo(k)fluoranthene	15.83	252	697718	37.07	nq	# 95
89) Benzo(a)pyrene	16.14	252	653187	36.87	nq	98
90) Dibenzo(a,h)anthracene	17.55	278	544210	36.28	nq	97
91) Benzo(g,h,i)perylene	17.93	276	558580	36.95	nq	95

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

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