

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070717\
 Data File : BF096555.D
 Acq On : 7 Jul 2017 13:26
 Operator : SJ/MA
 Sample : I4064-03MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PET-BF001-01MSD

Manual Integrations
 APPROVED

mohammad
 7/10/2017 2:22:13 PM

Quant Time: Jul 07 14:57:46 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 07 13:02:33 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	105999	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	453761	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	229828	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	400190	20.00	ng	0.00
75) Chrysene-d12	13.84	240	221018	20.00	ng	0.00
86) Perylene-d12	15.24	264	206360	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.32	112	905013	126.02	ng	0.01
7) Phenol-d6	6.35	99	1002259	113.52	ng	0.00
23) Nitrobenzene-d5	7.26	82	670470	88.79	ng	0.00
41) 2,4,6-Tribromophenol	10.52	330	214502	119.48	ng	0.00
44) 2-Fluorobiphenyl	9.06	172	1137030	84.59	ng	0.00
78) Terphenyl-d14	12.80	244	870760	84.19	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.40	88	128458	37.722	ng	# 76
3) Pyridine	3.10	79	315000	33.713	ng	# 67
4) n-Nitrosodimethylamine	3.04	42	210288	45.598	ng	# 81
6) Aniline	6.36	93	398790	34.948	ng	# 68
8) 2-Chlorophenol	6.49	128	317655	41.478	ng	# 95
9) Benzaldehyde	6.24	77	218523	39.548	ng	# 98
10) Phenol	6.36	94	404611	40.227	ng	# 76
11) bis(2-Chloroethyl)ether	6.44	93	301443	38.780	ng	# 93
12) 1,3-Dichlorobenzene	6.64	146	333626	41.547	ng	# 99
13) 1,4-Dichlorobenzene	6.72	146	340018	42.364	ng	# 95
14) 1,2-Dichlorobenzene	6.87	146	316456	42.942	ng	# 97
15) Benzyl Alcohol	6.85	79	264202	44.767	ng	# 98
16) 2,2'-oxybis(1-Chloropropan	6.98	45	654720	41.432	ng	# 91
17) 2-Methylphenol	6.96	107	262943	43.082	ng	# 97
18) Hexachloroethane	7.21	117	133256	45.796	ng	# 81
19) n-Nitroso-di-n-propylamine	7.12	70	230869	43.894	ng	# 85
20) 3+4-Methylphenols	7.12	107	336735	49.465	ng	# 87
22) Acetophenone	7.11	105	454708	45.862	ng	# 94
24) Nitrobenzene	7.28	77	361186	45.594	ng	# 93
25) Isophorone	7.53	82	676748	46.222	ng	# 95
26) 2-Nitrophenol	7.60	139	199969	47.692	ng	# 87
27) 2,4-Dimethylphenol	7.65	122	335816	47.037	ng	# 95
28) bis(2-Chloroethoxy)methane	7.73	93	409783	42.399	ng	# 99
29) 2,4-Dichlorophenol	7.85	162	275125	44.665	ng	# 99
30) 1,2,4-Trichlorobenzene	7.93	180	251930	43.357	ng	# 99
31) Naphthalene	8.00	128	930951	43.458	ng	# 100
32) Benzoic acid	7.76	122	116283	19.393	ng	# 90
33) 4-Chloroaniline	8.05	127	336887	37.862	ng	# 96
34) Hexachlorobutadiene	8.13	225	131420	44.097	ng	# 98
35) Caprolactam	8.42	113	108612m	51.133	ng	#
36) 4-Chloro-3-methylphenol	8.55	107	326695	47.932	ng	# 90
37) 2-Methylnaphthalene	8.70	142	692805	50.807	ng	# 98
39) 1,2,4,5-Tetrachlorobenzene	8.86	216	228808	38.342	ng	# 98
40) Hexachlorocyclopentadiene	8.85	237	224622	77.413	ng	# 99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.97	196	191172	40.496	ng	99
43) 2,4,5-Trichlorophenol	9.02	196	198619	42.552	ng	98
45) 1,1'-Biphenyl	9.16	154	769924	39.098	ng	98
46) 2-Chloronaphthalene	9.18	162	602182	41.034	ng	# 93
47) 2-Nitroaniline	9.27	65	226630	42.441	ng	96
48) Acenaphthylene	9.60	152	908581	39.072	ng	97
49) Dimethylphthalate	9.46	163	821349	47.873	ng	99
50) 2,6-Dinitrotoluene	9.52	165	165759	43.690	ng	# 87
51) Acenaphthene	9.77	154	556736	38.841	ng	98
52) 3-Nitroaniline	9.69	138	170172	35.961	ng	92
53) 2,4-Dinitrophenol	9.79	184	93949	46.622	ng	# 72
54) Dibenzofuran	9.94	168	787142	41.594	ng	98
55) 4-Nitrophenol	9.86	139	281720	71.827	ng	# 63
56) 2,4-Dinitrotoluene	9.93	165	203729	42.403	ng	# 90
57) Fluorene	10.28	166	560850	41.962	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.06	232	154471	47.833	ng	# 99
59) Diethylphthalate	10.16	149	691647	42.650	ng	99
60) 4-Chlorophenyl-phenylether	10.27	204	236380	40.649	ng	93
61) 4-Nitroaniline	10.30	138	176748	41.077	ng	91
62) Azobenzene	10.43	77	626553	39.957	ng	93
64) 4,6-Dinitro-2-methylphenol	10.33	198	83282	30.175	ng	82
65) n-Nitrosodiphenylamine	10.39	169	542221	38.616	ng	99
66) 4-Bromophenyl-phenylether	10.76	248	159232	40.874	ng	96
67) Hexachlorobenzene	10.83	284	166097	38.948	ng	# 94
68) Atrazine	10.92	200	169342	42.214	ng	94
69) Pentachlorophenol	11.02	266	166125	65.728	ng	98
70) Phenanthrene	11.24	178	905506	41.857	ng	100
71) Anthracene	11.29	178	928270	42.108	ng	99
72) Carbazole	11.44	167	907440	40.930	ng	98
73) Di-n-butylphthalate	11.78	149	1120500	41.727	ng	99
74) Fluoranthene	12.42	202	895846	42.237	ng	98
76) Benzidine	12.54	184	438093	41.310	ng	# 92
77) Pyrene	12.65	202	903832	50.947	ng	98
79) Butylbenzylphthalate	13.27	149	449690	50.074	ng	# 85
80) Benzo(a)anthracene	13.83	228	579513	45.279	ng	98
81) 3,3'-Dichlorobenzidine	13.79	252	175413	33.369	ng	# 95
82) Chrysene	13.87	228	595222	44.410	ng	99
83) Bis(2-ethylhexyl)phthalate	13.84	149	524158	52.288	ng	# 100
84) Di-n-octyl phthalate	14.44	149	930676	47.145	ng	# 95
85) Indeno(1,2,3-cd)pyrene	16.60	276	568521	53.737	ng	97
87) Benzo(b)fluoranthene	14.84	252	536422	41.485	ng	# 97
88) Benzo(k)fluoranthene	14.87	252	441198	38.145	ng	97
89) Benzo(a)pyrene	15.19	252	463292	40.132	ng	98
90) Dibenzo(a,h)anthracene	16.61	278	460766	50.735	ng	96
91) Benzo(g,h,i)perylene	17.00	276	503930	53.548	ng	97

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

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