

Data Path : Z:\HPCHEM1\BNA F\DATA\BF123115\
 Data File : BF084209.D
 Acq On : 31 Dec 2015 21:54
 Operator : UM/SJ
 Sample : PB87627BS
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB87627BS

Manual Integrations
 APPROVED

UMANGI
 1/4/2016 4:29:04 PM

Quant Time: Jan 02 00:05:07 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 31 14:15:57 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	277819	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	1182841	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	538685	20.00	ng	-0.01
63) Phenanthrene-d10	11.43	188	981545	20.00	ng	0.00
75) Chrysene-d12	14.07	240	730757	20.00	ng	0.00
86) Perylene-d12	15.51	264	535804	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1349201	78.55	ng	0.01
7) Phenol-d6	6.53	99	1907374	88.42	ng	0.00
23) Nitrobenzene-d5	7.47	82	1802183	84.80	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	443084	81.47	ng	-0.01
44) 2-Fluorobiphenyl	9.26	172	2522415	82.16	ng	0.00
78) Terphenyl-d14	13.01	244	2339134	71.45	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.61	88	306793	37.71	ng	# 71
3) Pyridine	3.34	79	925923	40.82	ng	# 74
4) n-Nitrosodimethylamine	3.30	42	490659	42.14	ng	78
6) Aniline	6.56	93	802016	25.41	ng	# 82
8) 2-Chlorophenol	6.68	128	814907	41.82	ng	87
9) Benzaldehyde	6.44	77	351283	25.01	ng	95
10) Phenol	6.54	94	1164711	47.49	ng	85
11) bis(2-Chloroethyl)ether	6.64	93	776708	40.88	ng	# 81
12) 1,3-Dichlorobenzene	6.84	146	865104m	43.03	ng	
13) 1,4-Dichlorobenzene	6.92	146	883196	43.81	ng	94
14) 1,2-Dichlorobenzene	7.07	146	772927	40.04	ng	97
15) Benzyl Alcohol	7.05	79	813509	44.83	ng	94
16) 2,2'-oxybis(1-Chloropropan	7.18	45	1418776	41.01	ng	84
17) 2-Methylphenol	7.15	107	709898	43.47	ng	96
18) Hexachloroethane	7.41	117	336954	41.80	ng	92
19) n-Nitroso-di-n-propylamine	7.32	70	638208	43.70	ng	# 72
20) 3+4-Methylphenols	7.31	107	842897	43.91	ng	# 80
22) Acetophenone	7.31	105	1106511	38.90	ng	# 82
24) Nitrobenzene	7.48	77	828358	38.43	ng	96
25) Isophorone	7.72	82	1677624	41.27	ng	97
26) 2-Nitrophenol	7.80	139	418372	42.65	ng	# 81
27) 2,4-Dimethylphenol	7.84	122	806938	40.64	ng	96
28) bis(2-Chloroethoxy)methane	7.94	93	1014828	40.87	ng	98
29) 2,4-Dichlorophenol	8.04	162	687809	41.58	ng	97
30) 1,2,4-Trichlorobenzene	8.13	180	732390	42.89	ng	# 95
31) Naphthalene	8.21	128	2408787	44.89	ng	99
32) Benzoic acid	7.95	122	486891	39.40	ng	91
33) 4-Chloroaniline	8.26	127	384847	15.64	ng	# 91
34) Hexachlorobutadiene	8.33	225	399161	40.66	ng	97
35) Caprolactam	8.62	113	224131	40.19	ng	# 37
36) 4-Chloro-3-methylphenol	8.74	107	788308	42.55	ng	96
37) 2-Methylnaphthalene	8.90	142	1488620	38.64	ng	96
39) 1,2,4,5-Tetrachlorobenzene	9.07	216	673428	43.49	ng	98
40) Hexachlorocyclopentadiene	9.06	237	864070	92.43	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.17	196	478277	41.28	nq	96
43) 2,4,5-Trichlorophenol	9.22	196	501234	45.13	nq #	93
45) 1,1'-Biphenyl	9.37	154	1904954	44.49	nq	100
46) 2-Chloronaphthalene	9.39	162	1407586	41.86	nq	95
47) 2-Nitroaniline	9.48	65	476862	42.96	nq	97
48) Acenaphthylene	9.80	152	2109326	42.46	nq	97
49) Dimethylphthalate	9.68	163	1684792	43.75	nq #	91
50) 2,6-Dinitrotoluene	9.73	165	374493	44.34	nq	92
51) Acenaphthene	9.97	154	1551088	46.54	nq	97
52) 3-Nitroaniline	9.89	138	250206	23.91	nq #	87
53) 2,4-Dinitrophenol	10.00	184	285613	79.41	nq #	77
54) Dibenzofuran	10.16	168	2007174	46.61	nq	98
55) 4-Nitrophenol	10.05	139	603925	79.49	nq #	78
56) 2,4-Dinitrotoluene	10.13	165	553836	51.81	nq #	84
57) Fluorene	10.49	166	1417279	42.32	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.27	232	417762	44.05	nq	89
59) Diethylphthalate	10.37	149	1758539	46.31	nq	96
60) 4-Chlorophenyl-phenylether	10.49	204	719933	52.43	nq	99
61) 4-Nitroaniline	10.51	138	409329	43.87	nq	93
62) Azobenzene	10.65	77	1937349	46.52	nq	93
64) 4,6-Dinitro-2-methylphenol	10.53	198	228681	38.15	nq #	68
65) n-Nitrosodiphenylamine	10.60	169	1251359	39.87	nq	99
66) 4-Bromophenyl-phenylether	10.98	248	483459	45.76	nq	95
67) Hexachlorobenzene	11.04	284	467623	39.33	nq #	79
68) Atrazine	11.14	200	470001	45.76	nq	90
69) Pentachlorophenol	11.23	266	519059	84.19	nq	99
70) Phenanthrene	11.45	178	2246930	43.76	nq	96
71) Anthracene	11.51	178	2253207	44.77	nq	96
72) Carbazole	11.65	167	1918168	38.98	nq	98
73) Di-n-butylphthalate	12.00	149	2551413	50.23	nq #	95
74) Fluoranthene	12.64	202	2117473	39.48	nq	97
76) Benzidine	12.76	184	700300	26.73	nq	99
77) Pyrene	12.87	202	2149971	37.49	nq	97
79) Butylbenzylphthalate	13.49	149	1108413	44.67	nq #	97
80) Benzo(a)anthracene	14.05	228	1657582	38.63	nq	99
81) 3,3'-Dichlorobenzidine	14.02	252	371100	24.78	nq #	97
82) Chrysene	14.09	228	1535809	37.25	nq	98
83) Bis(2-ethylhexyl)phthalate	14.05	149	1299286	41.44	nq #	96
84) Di-n-octyl phthalate	14.66	149	2034289	38.44	nq #	86
85) Indeno(1,2,3-cd)pyrene	16.92	276	1403695	42.95	nq	99
87) Benzo(b)fluoranthene	15.08	252	1659646m	47.35	nq	
88) Benzo(k)fluoranthene	15.12	252	1069614m	37.31	nq	
89) Benzo(a)pyrene	15.44	252	1301294	43.77	nq #	95
90) Dibenzo(a,h)anthracene	16.95	278	1144339	44.73	nq	95
91) Benzo(g,h,i)perylene	17.35	276	1191956	44.99	nq	97

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

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