

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080415\
 Data File : BF080879.D
 Acq On : 5 Aug 2015 13:13
 Operator : TP/UM
 Sample : G3188-01MSD
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-EF-01-003-ABCDMSD

Manual Integrations
 APPROVED

MMDadoda
 8/5/2015 7:02:05 PM

Quant Time: Aug 05 16:23:40 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 05 01:46:16 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	49147	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	176575	20.00	ng	0.00
38) Acenaphthene-d10	9.86	164	68933	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	103488	20.00	ng	0.00
75) Chrysene-d12	13.97	240	94964	20.00	ng	0.01
86) Perylene-d12	15.38	264	124976	20.00	ng	0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	412953	127.17	ng	0.01
7) Phenol-d6	6.46	99	528593	124.52	ng	0.01
23) Nitrobenzene-d5	7.39	82	320099	85.43	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	83498	125.61	ng	0.00
44) 2-Fluorobiphenyl	9.18	172	465490	90.93	ng	0.00
78) Terphenyl-d14	12.92	244	303765	57.89	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	2.50	88	55978	35.16	ng	98
3) Pyridine	3.21	79	151242	34.47	ng	98
4) n-Nitrosodimethylamine	3.15	42	91841	40.81	ng	97
6) Aniline	6.49	93	172222	27.88	ng	100
8) 2-Chlorophenol	6.61	128	146523	40.28	ng	95
9) Benzaldehyde	6.37	77	27800	9.39	ng	96
10) Phenol	6.48	94	210165	41.61	ng	99
11) bis(2-Chloroethyl)ether	6.57	93	161230	42.75	ng	98
12) 1,3-Dichlorobenzene	6.76	146	140949	36.16	ng	99
13) 1,4-Dichlorobenzene	6.84	146	143122	36.58	ng	98
14) 1,2-Dichlorobenzene	7.00	146	137467	37.24	ng	99
15) Benzyl Alcohol	6.97	79	121900	43.97	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.10	45	295724	40.51	ng	99
17) 2-Methylphenol	7.08	107	120830	37.92	ng	98
18) Hexachloroethane	7.34	117	53424	35.93	ng	96
19) n-Nitroso-di-n-propylamine	7.24	70	111974	40.83	ng	96
20) 3+4-Methylphenols	7.23	107	137550	37.23	ng	# 79
22) Acetophenone	7.24	105	193119	44.78	ng	# 96
24) Nitrobenzene	7.40	77	149603	38.50	ng	# 83
25) Isophorone	7.65	82	282226	42.51	ng	97
26) 2-Nitrophenol	7.72	139	62839	39.42	ng	100
27) 2,4-Dimethylphenol	7.77	122	128060	43.15	ng	96
28) bis(2-Chloroethoxy)methane	7.86	93	167766	41.09	ng	99
29) 2,4-Dichlorophenol	7.96	162	95006	38.77	ng	99
30) 1,2,4-Trichlorobenzene	8.05	180	112858	40.48	ng	98
31) Naphthalene	8.13	128	393947	40.96	ng	100
32) Benzoic acid	7.82	122	10912	7.78	ng	96
33) 4-Chloroaniline	8.18	127	101322	26.93	ng	96
34) Hexachlorobutadiene	8.25	225	61930	38.96	ng	97
35) Caprolactam	8.53	113	30484	43.57	ng	# 87
36) 4-Chloro-3-methylphenol	8.66	107	109370	40.68	ng	95
37) 2-Methylnaphthalene	8.82	142	234748	40.73	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.99	216	92900	39.49	ng	99
40) Hexachlorocyclopentadiene	8.98	237	64507	45.66	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.09	196	59696	37.44	ng	98
43) 2,4,5-Trichlorophenol	9.14	196	64635	40.42	ng	# 84
45) 1,1'-Biphenyl	9.29	154	263015	41.39	ng	99
46) 2-Chloronaphthalene	9.31	162	197930	40.05	ng	98
47) 2-Nitroaniline	9.40	65	81184	45.12	ng	97
48) Acenaphthylene	9.72	152	340720	41.90	ng	100
49) Dimethylphthalate	9.58	163	312448	62.15	ng	100
50) 2,6-Dinitrotoluene	9.64	165	43890	41.74	ng	97
51) Acenaphthene	9.89	154	194574	39.80	ng	100
52) 3-Nitroaniline	9.81	138	41230	31.75	ng	92
53) 2,4-Dinitrophenol	9.92	184	13025	24.76	ng	93
54) Dibenzofuran	10.06	168	284698	43.67	ng	99
55) 4-Nitrophenol	9.97	139	75301	82.42	ng	# 75
56) 2,4-Dinitrotoluene	10.04	165	59308	44.68	ng	99
57) Fluorene	10.41	166	199064	39.98	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.18	232	44349	38.64	ng	97
59) Diethylphthalate	10.28	149	229677	46.15	ng	99
60) 4-Chlorophenyl-phenylether	10.40	204	89160	43.82	ng	98
61) 4-Nitroaniline	10.42	138	43946	37.92	ng	93
62) Azobenzene	10.56	77	231564	39.78	ng	99
64) 4,6-Dinitro-2-methylphenol	10.44	198	11916	17.90	ng	# 1
65) n-Nitrosodiphenylamine	10.51	169	162920	41.09	ng	97
66) 4-Bromophenyl-phenylether	10.89	248	52961	46.24	ng	97
67) Hexachlorobenzene	10.96	284	50218	40.14	ng	# 61
68) Atrazine	11.05	200	50425	48.86	ng	98
69) Pentachlorophenol	11.14	266	48687	71.42	ng	98
70) Phenanthrene	11.36	178	244392	40.87	ng	99
71) Anthracene	11.41	178	266925	45.94	ng	98
72) Carbazole	11.56	167	218838	39.93	ng	99
73) Di-n-butylphthalate	11.90	149	317480	48.92	ng	100
74) Fluoranthene	12.55	202	266962	45.84	ng	99
76) Benzidine	12.67	184	192922	41.97	ng	99
77) Pyrene	12.77	202	278684	32.71	ng	99
79) Butylbenzylphthalate	13.40	149	116153	33.34	ng	# 65
80) Benzo(a)anthracene	13.96	228	284369	44.77	ng	98
81) 3,3'-Dichlorobenzidine	13.93	252	105220	44.58	ng	# 94
82) Chrysene	14.00	228	280869	43.39	ng	98
83) Bis(2-ethylhexyl)phthalate	13.96	149	166546	35.80	ng	98
84) Di-n-octyl phthalate	14.57	149	336565	43.04	ng	97
85) Indeno(1,2,3-cd)pyrene	16.75	276	317110	59.67	ng	99
87) Benzo(b)fluoranthene	14.98	252	343470m	36.74	ng	
88) Benzo(k)fluoranthene	15.01	252	321487m	44.68	ng	
89) Benzo(a)pyrene	15.32	252	323921	42.45	ng	98
90) Dibenzo(a,h)anthracene	16.76	278	263000	39.67	ng	97
91) Benzo(g,h,i)perylene	17.16	276	251431	39.24	ng	97

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

