

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071415\
 Data File : BF080488.D
 Acq On : 15 Jul 2015 13:00
 Operator : TP/UM
 Sample : G2973-10MSD
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-7D78MSD

Manual Integrations
 APPROVED

apatel
 7/16/2015 8:21:55 AM

Quant Time: Jul 15 16:08:04 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 15 03:58:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.13	152	23756	20.00	ng	0.00
21) Naphthalene-d8	8.41	136	102468	20.00	ng	0.00
38) Acenaphthene-d10	10.17	164	51104	20.00	ng	-0.01
63) Phenanthrene-d10	11.66	188	111839	20.00	ng	0.00
75) Chrysene-d12	14.67	240	129115	20.00	ng	0.19
86) Perylene-d12	16.56	264	116806	20.00	ng	0.33

System Monitoring Compounds

5) 2-Fluorophenol	5.76	112	139510	95.99	ng	0.00
7) Phenol-d6	6.76	99	190206	106.94	ng	-0.01
23) Nitrobenzene-d5	7.69	82	120607	64.43	ng	0.00
41) 2,4,6-Tribromophenol	10.97	330	49332	102.13	ng	0.00
44) 2-Fluorobiphenyl	9.48	172	216769	58.12	ng	0.00
78) Terphenyl-d14	13.37	244	280159	52.30	ng	0.09

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.86	88	17569	30.01	ng	# 91
3) Pyridine	3.84	79	34378	26.57	ng	96
4) n-Nitrosodimethylamine	3.66	42	26635	35.02	ng	97
6) Aniline	6.80	93	61211	23.71	ng	99
8) 2-Chlorophenol	6.92	128	53506	35.09	ng	94
9) Benzaldehyde	6.68	77	20499	18.87	ng	98
10) Phenol	6.78	94	71155	34.70	ng	96
11) bis(2-Chloroethyl)ether	6.85	93	54754	35.13	ng	100
12) 1,3-Dichlorobenzene	7.06	146	58460m	32.88	ng	
13) 1,4-Dichlorobenzene	7.14	146	64640	33.44	ng	97
14) 1,2-Dichlorobenzene	7.30	146	61791	33.07	ng	96
15) Benzyl Alcohol	7.28	79	43375	34.08	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.39	45	91351	36.16	ng	78
17) 2-Methylphenol	7.39	107	44737	36.25	ng	# 90
18) Hexachloroethane	7.64	117	20967	32.92	ng	93
19) n-Nitroso-di-n-propylamine	7.53	70	39743	32.99	ng	99
20) 3+4-Methylphenols	7.54	107	63186	36.13	ng	# 88
22) Acetophenone	7.53	105	78282	32.56	ng	# 96
24) Nitrobenzene	7.71	77	62186	30.14	ng	95
25) Isophorone	7.94	82	118499	34.28	ng	96
26) 2-Nitrophenol	8.03	139	27335	30.54	ng	94
27) 2,4-Dimethylphenol	8.06	122	53381	32.52	ng	99
28) bis(2-Chloroethoxy)methane	8.14	93	66928	34.43	ng	99
29) 2,4-Dichlorophenol	8.27	162	49148	34.34	ng	91
30) 1,2,4-Trichlorobenzene	8.35	180	53104	29.82	ng	97
31) Naphthalene	8.43	128	169221	31.93	ng	100
32) Benzoic acid	8.14	122	20298	29.81	ng	85
33) 4-Chloroaniline	8.49	127	48508	22.99	ng	98
34) Hexachlorobutadiene	8.54	225	27883	30.90	ng	99
35) Caprolactam	8.83	113	12685	34.35	ng	# 69
36) 4-Chloro-3-methylphenol	8.97	107	50100	35.76	ng	# 76
37) 2-Methylnaphthalene	9.13	142	113294	31.38	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.29	216	48992	28.89	ng	98
40) Hexachlorocyclopentadiene	9.28	237	41851	53.16	ng	94

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42) 2,4,6-Trichlorophenol	9.41	196	31670	30.04	ng	98
43) 2,4,5-Trichlorophenol	9.45	196	34135	32.14	ng #	75
45) 1,1'-Biphenyl	9.58	154	143909	32.72	ng	98
46) 2-Chloronaphthalene	9.62	162	108015	29.90	ng	97
47) 2-Nitroaniline	9.72	65	33931	32.56	ng	92
48) Acenaphthylene	10.03	152	173533	31.76	ng	99
49) Dimethylphthalate	9.88	163	189114	46.75	ng	99
50) 2,6-Dinitrotoluene	9.95	165	29559	33.95	ng	93
51) Acenaphthene	10.20	154	111131	33.44	ng	96
52) 3-Nitroaniline	10.13	138	25497	27.21	ng	94
53) 2,4-Dinitrophenol	10.24	184	19637	58.13	ng	90
54) Dibenzofuran	10.37	168	151923	32.41	ng	99
55) 4-Nitrophenol	10.30	139	43443	63.80	ng #	85
56) 2,4-Dinitrotoluene	10.35	165	34680	35.16	ng #	23
57) Fluorene	10.72	166	123540	32.78	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.50	232	29821	34.67	ng	99
59) Diethylphthalate	10.58	149	135185	33.77	ng	100
60) 4-Chlorophenyl-phenylether	10.70	204	60680	34.33	ng	93
61) 4-Nitroaniline	10.75	138	27237	32.03	ng	94
62) Azobenzene	10.86	77	137341	35.90	ng	98
64) 4,6-Dinitro-2-methylphenol	10.76	198	15860	29.93	ng	86
65) n-Nitrosodiphenylamine	10.83	169	106876	29.42	ng	99
66) 4-Bromophenyl-phenylether	11.20	248	35233	31.59	ng	88
67) Hexachlorobenzene	11.28	284	36774	30.30	ng	98
68) Atrazine	11.34	200	35293	33.35	ng	98
69) Pentachlorophenol	11.47	266	38367	68.62	ng	96
70) Phenanthrene	11.69	178	194467	29.92	ng	98
71) Anthracene	11.74	178	212597	34.79	ng	99
72) Carbazole	11.90	167	183086	33.07	ng	98
73) Di-n-butylphthalate	12.22	149	223251	38.00	ng #	98
74) Fluoranthene	12.96	202	234136	36.53	ng	98
76) Benzidine	13.09	184	121235	31.69	ng	99
77) Pyrene	13.21	202	240900	31.69	ng	100
79) Butylbenzylphthalate	13.93	149	101487	36.55	ng	95
80) Benzo(a)anthracene	14.66	228	233745	31.06	ng	100
81) 3,3'-Dichlorobenzidine	14.61	252	67708	28.47	ng #	95
82) Chrysene	14.71	228	220000	30.58	ng	98
83) Bis(2-ethylhexyl)phthalate	14.63	149	155074	38.45	ng #	95
84) Di-n-octyl phthalate	15.47	149	254890	35.19	ng	99
85) Indeno(1,2,3-cd)pyrene	18.23	276	233137	23.79	ng	99
87) Benzo(b)fluoranthene	16.05	252	217099	32.53	ng #	98
88) Benzo(k)fluoranthene	16.09	252	222559	34.52	ng	100
89) Benzo(a)pyrene	16.48	252	210924	32.09	ng	98
90) Dibenzo(a,h)anthracene	18.24	278	192982	25.89	ng	98
91) Benzo(g,h,i)perylene	18.73	276	193755	24.84	ng	99

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

