

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
 Data File : BF087486.D
 Acq On : 28 May 2016 18:51
 Operator : UM/SJ
 Sample : H3258-02MS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-2MS

Manual Integrations
 APPROVED

umangi
 5/31/2016 7:03:49 PM

Quant Time: May 30 03:37:16 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 30 03:25:10 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	110323	20.00	ng	-0.01
21) Naphthalene-d8	9.50	136	457497	20.00	ng	-0.01
38) Acenaphthene-d10	12.33	164	215173	20.00	ng	0.00
63) Phenanthrene-d10	14.73	188	397603	20.00	ng	0.00
75) Chrysene-d12	18.43	240	316330	20.00	ng	0.00
86) Perylene-d12	20.10	264	216331	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.37	112	784838	125.00	ng	0.01
7) Phenol-d6	7.03	99	1029265	121.32	ng	-0.01
23) Nitrobenzene-d5	8.39	82	709913	78.21	ng	-0.01
41) 2,4,6-Tribromophenol	13.62	330	235816	114.04	ng	-0.01
44) 2-Fluorobiphenyl	11.28	172	1175414	77.01	ng	-0.01
78) Terphenyl-d14	17.09	244	1102215	74.08	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.83	88	91791	27.71	ng	# 65
3) Pyridine	2.41	79	204269	28.20	ng	# 64
4) n-Nitrosodimethylamine	2.38	42	203517	36.47	ng	# 46
6) Aniline	6.97	93	321507	29.29	ng	94
8) 2-Chlorophenol	7.15	128	309316	39.50	ng	95
9) Benzaldehyde	6.81	77	45918	9.80	ng	94
10) Phenol	7.04	94	392589	42.38	ng	76
11) bis(2-Chloroethyl)ether	7.11	93	288118	38.43	ng	84
12) 1,3-Dichlorobenzene	7.36	146	312502	36.59	ng	# 91
13) 1,4-Dichlorobenzene	7.50	146	323656	36.92	ng	97
14) 1,2-Dichlorobenzene	7.72	146	311156	38.37	ng	99
15) Benzyl Alcohol	7.77	79	266928	43.16	ng	90
16) 2,2'-oxybis(1-Chloropropan	7.96	45	594255	35.22	ng	87
17) 2-Methylphenol	7.99	107	256191	40.83	ng	# 89
18) Hexachloroethane	8.24	117	122780	37.54	ng	94
19) n-Nitroso-di-n-propylamine	8.19	70	256500	40.54	ng	# 78
20) 3+4-Methylphenols	8.25	107	333731	44.00	ng	# 61
22) Acetophenone	8.16	105	452921	41.44	ng	# 97
24) Nitrobenzene	8.41	77	360059	37.58	ng	96
25) Isophorone	8.81	82	630352	38.72	ng	# 97
26) 2-Nitrophenol	8.92	139	159704	40.77	ng	# 80
27) 2,4-Dimethylphenol	9.06	122	262205	38.56	ng	# 84
28) bis(2-Chloroethoxy)methane	9.19	93	369103	40.25	ng	98
29) 2,4-Dichlorophenol	9.34	162	259513	42.35	ng	97
30) 1,2,4-Trichlorobenzene	9.43	180	267981	38.20	ng	98
31) Naphthalene	9.53	128	889855	38.82	ng	100
32) Benzoic acid	9.34	122	11914	8.81	ng	# 73
33) 4-Chloroaniline	9.68	127	221653	26.25	ng	# 89
34) Hexachlorobutadiene	9.75	225	156534	36.71	ng	97
35) Caprolactam	10.30	113	71283m	39.86	ng	
36) 4-Chloro-3-methylphenol	10.55	107	286094	41.30	ng	91
37) 2-Methylnaphthalene	10.66	142	581532	40.61	ng	98
39) 1,2,4,5-Tetrachlorobenzene	10.94	216	256381	37.22	ng	99
40) Hexachlorocyclopentadiene	10.90	237	145159	54.62	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	11.16	196	149664	37.65	ng	91
43) 2,4,5-Trichlorophenol	11.24	196	147453	36.42	ng #	86
45) 1,1'-Biphenyl	11.43	154	715841	39.51	ng	98
46) 2-Chloronaphthalene	11.44	162	524985	37.88	ng #	92
47) 2-Nitroaniline	11.67	65	218015	43.66	ng	85
48) Acenaphthylene	12.10	152	854200	38.93	ng	99
49) Dimethylphthalate	11.99	163	892144	51.76	ng	100
50) 2,6-Dinitrotoluene	12.08	165	136960	39.43	ng #	82
51) Acenaphthene	12.38	154	508025	37.67	ng	99
52) 3-Nitroaniline	12.34	138	109814	30.81	ng #	68
53) 2,4-Dinitrophenol	12.57	184	59048	37.05	ng #	81
54) Dibenzofuran	12.66	168	762885	41.78	ng	93
55) 4-Nitrophenol	12.74	139	128904	75.81	ng #	39
56) 2,4-Dinitrotoluene	12.74	165	187657	40.43	ng #	91
57) Fluorene	13.22	166	612722	38.70	ng	100
58) 2,3,4,6-Tetrachlorophenol	12.91	232	133668	42.92	ng	96
59) Diethylphthalate	13.13	149	626241	37.37	ng	100
60) 4-Chlorophenyl-phenylether	13.25	204	282438	36.58	ng	96
61) 4-Nitroaniline	13.35	138	131773	39.61	ng #	63
62) Azobenzene	13.51	77	762098	43.84	ng	86
64) 4,6-Dinitro-2-methylphenol	13.41	198	81959	32.42	ng #	37
65) n-Nitrosodiphenylamine	13.46	169	510598	39.71	ng	99
66) 4-Bromophenyl-phenylether	14.03	248	165825	38.12	ng	91
67) Hexachlorobenzene	14.10	284	170769	37.54	ng #	73
68) Atrazine	14.39	200	167965	40.98	ng	93
69) Pentachlorophenol	14.47	266	81257	59.21	ng	99
70) Phenanthrene	14.77	178	870001	39.50	ng	99
71) Anthracene	14.85	178	875310	39.34	ng	99
72) Carbazole	15.14	167	783607	41.30	ng	99
73) Di-n-butylphthalate	15.73	149	959389	39.10	ng	98
74) Fluoranthene	16.54	202	893703	40.47	ng	91
76) Benzidine	16.77	184	211659	26.00	ng	95
77) Pyrene	16.83	202	874819	38.42	ng	99
79) Butylbenzylphthalate	17.75	149	383411	37.97	ng #	74
80) Benzo(a)anthracene	18.42	228	711779	39.56	ng	99
81) 3,3'-Dichlorobenzidine	18.41	252	130006	13.08	ng #	96
82) Chrysene	18.47	228	677740	37.95	ng	99
83) Bis(2-ethylhexyl)phthalate	18.49	149	517565	35.13	ng #	95
84) Di-n-octyl phthalate	19.27	149	777631	34.61	ng #	86
85) Indeno(1,2,3-cd)pyrene	21.31	276	508754	35.54	ng	94
87) Benzo(b)fluoranthene	19.69	252	586490	42.56	ng	97
88) Benzo(k)fluoranthene	19.73	252	446839m	37.67	ng	
89) Benzo(a)pyrene	20.05	252	478466	40.15	ng	99
90) Dibenzo(a,h)anthracene	21.31	278	429556	42.26	ng	96
91) Benzo(g,h,i)perylene	21.67	276	431572	43.03	ng #	91

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

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