

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
 Data File : BF085699.D
 Acq On : 23 Mar 2016 11:19
 Operator : UM/SJ
 Sample : H1857-18MSD
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-06(8-10)MSD

Manual Integrations
 APPROVED

UMANGI
 3/24/2016 11:49:16 AM

Quant Time: Mar 23 14:55:46 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 18 23:31:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	127162	20.00	ng	-0.02
21) Naphthalene-d8	8.12	136	499236	20.00	ng	-0.02
38) Acenaphthene-d10	9.88	164	239949	20.00	ng	-0.02
63) Phenanthrene-d10	11.35	188	369692	20.00	ng	-0.03
75) Chrysene-d12	13.99	240	233068	20.00	ng	-0.02
86) Perylene-d12	15.41	264	217663	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.44	112	1002048	132.36	ng	-0.01
7) Phenol-d6	6.48	99	1349720	139.01	ng	-0.01
23) Nitrobenzene-d5	7.41	82	909707	105.77	ng	-0.02
41) 2,4,6-Tribromophenol	10.66	330	282209	120.07	ng	-0.02
44) 2-Fluorobiphenyl	9.20	172	1310524	96.40	ng	-0.02
78) Terphenyl-d14	12.94	244	863066	89.11	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.48	88	130892	35.57	ng	97
3) Pyridine	3.20	79	325627	36.03	ng	99
4) n-Nitrosodimethylamine	3.14	42	154029	40.83	ng	96
6) Aniline	6.49	93	145301	11.14	ng	# 1
8) 2-Chlorophenol	6.62	128	363936	41.55	ng	91
9) Benzaldehyde	6.38	77	59932	10.11	ng	98
10) Phenol	6.49	94	528272	47.77	ng	98
11) bis(2-Chloroethyl)ether	6.57	93	537694	64.49	ng	98
12) 1,3-Dichlorobenzene	6.78	146	392206	41.09	ng	99
13) 1,4-Dichlorobenzene	6.85	146	380651	39.10	ng	100
14) 1,2-Dichlorobenzene	7.01	146	386287	44.69	ng	99
15) Benzyl Alcohol	6.98	79	321693	48.37	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.11	45	459393	41.90	ng	84
17) 2-Methylphenol	7.10	107	332525	45.80	ng	99
18) Hexachloroethane	7.35	117	131633	37.45	ng	93
19) n-Nitroso-di-n-propylamine	7.26	70	275776	47.55	ng	98
20) 3+4-Methylphenols	7.25	107	388195	46.17	ng	# 85
22) Acetophenone	7.25	105	503430	42.49	ng	97
24) Nitrobenzene	7.42	77	366686	38.95	ng	99
25) Isophorone	7.66	82	751602	43.65	ng	99
26) 2-Nitrophenol	7.74	139	215857	46.39	ng	97
27) 2,4-Dimethylphenol	7.78	122	394578	47.96	ng	94
28) bis(2-Chloroethoxy)methane	7.88	93	500864	51.14	ng	99
29) 2,4-Dichlorophenol	7.98	162	303811	44.71	ng	95
30) 1,2,4-Trichlorobenzene	8.06	180	309716	40.69	ng	98
31) Naphthalene	8.14	128	1092826	43.37	ng	100
32) Benzoic acid	7.78	122	394578	57.36	ng	# 11
33) 4-Chloroaniline	8.21	127	216911	21.00	ng	97
34) Hexachlorobutadiene	8.26	225	163706	40.36	ng	98
35) Caprolactam	8.57	113	109926m	47.02	ng	
36) 4-Chloro-3-methylphenol	8.69	107	368565	46.44	ng	91
37) 2-Methylnaphthalene	8.84	142	758810	49.40	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.00	216	265610	36.49	ng	99
40) Hexachlorocyclopentadiene	8.98	237	48065	14.65	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.11	196	203520	40.85	ng	99
43) 2,4,5-Trichlorophenol	9.16	196	210158	41.45	ng #	81
45) 1,1'-Biphenyl	9.29	154	776957	37.53	ng	99
46) 2-Chloronaphthalene	9.33	162	630473	41.70	ng	98
47) 2-Nitroaniline	9.42	65	195601	42.76	ng	95
48) Acenaphthylene	9.74	152	1014147	41.45	ng	99
49) Dimethylphthalate	9.60	163	945606	52.22	ng	100
50) 2,6-Dinitrotoluene	9.67	165	163493	42.77	ng #	52
51) Acenaphthene	9.91	154	633341	41.25	ng	100
52) 3-Nitroaniline	9.83	138	136554	28.52	ng	98
53) 2,4-Dinitrophenol	9.94	184	8765	3.74	ng #	80
54) Dibenzofuran	10.08	168	908216	48.57	ng	99
55) 4-Nitrophenol	10.00	139	259534	70.14	ng	98
56) 2,4-Dinitrotoluene	10.07	165	227774	45.51	ng #	77
57) Fluorene	10.42	166	674969	43.31	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.20	232	149461	40.94	ng	99
59) Diethylphthalate	10.30	149	754422	41.91	ng	99
60) 4-Chlorophenyl-phenylether	10.41	204	313087	41.14	ng	93
61) 4-Nitroaniline	10.45	138	187152	39.66	ng	98
62) Azobenzene	10.57	77	783350	45.35	ng	97
64) 4,6-Dinitro-2-methylphenol	10.48	198	29719	12.56	ng #	55
65) n-Nitrosodiphenylamine	10.54	169	587993	51.06	ng	98
66) 4-Bromophenyl-phenylether	10.90	248	188920	51.15	ng	92
67) Hexachlorobenzene	10.97	284	189317	48.25	ng	95
68) Atrazine	11.06	200	188056	53.64	ng	99
69) Pentachlorophenol	11.17	266	182379	76.56	ng	99
70) Phenanthrene	11.38	178	997443	51.17	ng	100
71) Anthracene	11.43	178	818668	40.05	ng	99
72) Carbazole	11.59	167	816447	46.30	ng	99
73) Di-n-butylphthalate	11.92	149	1028806	46.59	ng	100
74) Fluoranthene	12.57	202	760925	37.28	ng	99
76) Benzidine	12.69	184	268405	34.25	ng	99
77) Pyrene	12.80	202	805982	48.16	ng	100
79) Butylbenzylphthalate	13.41	149	337412	42.30	ng #	77
80) Benzo(a)anthracene	13.98	228	625188	46.21	ng	99
81) 3,3'-Dichlorobenzidine	13.95	252	155442	31.35	ng #	97
82) Chrysene	14.01	228	464174	35.66	ng	99
83) Bis(2-ethylhexyl)phthalate	13.97	149	458861	46.30	ng #	98
84) Di-n-octyl phthalate	14.57	149	764457	47.34	ng	99
85) Indeno(1,2,3-cd)pyrene	16.79	276	498480	45.83	ng	99
87) Benzo(b)fluoranthene	15.01	252	644523m	45.38	ng	
88) Benzo(k)fluoranthene	15.03	252	448234m	38.31	ng	
89) Benzo(a)pyrene	15.35	252	520438	43.22	ng	98
90) Dibenzo(a,h)anthracene	16.80	278	430760	42.89	ng #	96
91) Benzo(g,h,i)perylene	17.21	276	385766	39.69	ng	96

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

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