

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042016\
 Data File : BF086307.D
 Acq On : 20 Apr 2016 14:13
 Operator : UM/SJ
 Sample : SSTDICV040
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF042016

Manual Integrations
 APPROVED

sohil
 4/21/2016 7:09:03 PM

Quant Time: Apr 20 15:07:58 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 20 15:02:43 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	260947	20.00	ng	-0.03
21) Naphthalene-d8	8.16	136	1076421	20.00	ng	-0.02
38) Acenaphthene-d10	9.90	164	468383	20.00	ng	-0.03
63) Phenanthrene-d10	11.39	188	755215	20.00	ng	-0.02
75) Chrysene-d12	14.02	240	427455	20.00	ng	-0.03
86) Perylene-d12	15.45	264	470825	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.45	112	1209321	80.86	ng	-0.03
7) Phenol-d6	6.50	99	1497449	76.51	ng	-0.02
23) Nitrobenzene-d5	7.44	82	1432692	76.23	ng	-0.02
41) 2,4,6-Tribromophenol	10.69	330	292040	80.00	ng	-0.03
44) 2-Fluorobiphenyl	9.23	172	2026185	82.04	ng	-0.03
78) Terphenyl-d14	12.97	244	1337188	74.46	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.46	88	317843	39.23	ng	98
3) Pyridine	3.20	79	839694	40.48	ng	95
4) n-Nitrosodimethylamine	3.14	42	291986	40.86	ng	# 69
6) Aniline	6.52	93	1036199	39.34	ng	91
8) 2-Chlorophenol	6.65	128	667178	37.55	ng	89
9) Benzaldehyde	6.41	77	551594	39.76	ng	92
10) Phenol	6.51	94	920022	41.02	ng	100
11) bis(2-Chloroethyl)ether	6.60	93	746213	38.09	ng	88
12) 1,3-Dichlorobenzene	6.81	146	761936m	39.09	ng	
13) 1,4-Dichlorobenzene	6.89	146	779802	37.44	ng	96
14) 1,2-Dichlorobenzene	7.04	146	695964	37.67	ng	96
15) Benzyl Alcohol	7.01	79	483690	39.78	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.15	45	1097849	39.83	ng	92
17) 2-Methylphenol	7.12	107	572927	38.59	ng	98
18) Hexachloroethane	7.38	117	251767	40.56	ng	94
19) n-Nitroso-di-n-propylamine	7.29	70	472336	39.84	ng	97
20) 3+4-Methylphenols	7.28	107	594495	36.91	ng	# 79
22) Acetophenone	7.28	105	816349	36.10	ng	# 87
24) Nitrobenzene	7.45	77	744092	37.01	ng	94
25) Isophorone	7.69	82	1312668	36.86	ng	99
26) 2-Nitrophenol	7.77	139	385065	42.81	ng	# 79
27) 2,4-Dimethylphenol	7.80	122	643134	36.46	ng	95
28) bis(2-Chloroethoxy)methane	7.90	93	822453	39.49	ng	99
29) 2,4-Dichlorophenol	8.01	162	553914	39.82	ng	97
30) 1,2,4-Trichlorobenzene	8.10	180	563495	37.55	ng	96
31) Naphthalene	8.18	128	2016966	39.31	ng	97
32) Benzoic acid	7.93	122	424451	40.49	ng	93
33) 4-Chloroaniline	8.22	127	845177	39.80	ng	95
34) Hexachlorobutadiene	8.29	225	275262	36.40	ng	99
35) Caprolactam	8.60	113	186810	41.47	ng	# 78
36) 4-Chloro-3-methylphenol	8.70	107	617900	39.50	ng	98
37) 2-Methylnaphthalene	8.86	142	1169112	36.90	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.04	216	477078	38.27	ng	97
40) Hexachlorocyclopentadiene	9.01	237	111889	37.58	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.14	196	335713	38.82	nq	97
43) 2,4,5-Trichlorophenol	9.18	196	370433	39.33	nq	96
45) 1,1'-Biphenyl	9.33	154	1433886	38.98	nq	97
46) 2-Chloronaphthalene	9.36	162	1130839	40.57	nq	95
47) 2-Nitroaniline	9.45	65	381712	40.29	nq	88
48) Acenaphthylene	9.77	152	1752915	40.27	nq	100
49) Dimethylphthalate	9.63	163	1319780	37.01	nq	100
50) 2,6-Dinitrotoluene	9.69	165	288536	39.32	nq	# 67
51) Acenaphthene	9.94	154	1005525	39.38	nq	100
52) 3-Nitroaniline	9.86	138	369929	39.64	nq	92
53) 2,4-Dinitrophenol	9.96	184	52928	38.32	nq	# 65
54) Dibenzofuran	10.11	168	1389071	40.01	nq	96
55) 4-Nitrophenol	10.02	139	229953	39.74	nq	# 75
56) 2,4-Dinitrotoluene	10.09	165	357024	39.15	nq	# 68
57) Fluorene	10.45	166	1041646	40.49	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.22	232	245153	38.79	nq	99
59) Diethylphthalate	10.33	149	1270704	36.52	nq	99
60) 4-Chlorophenyl-phenylether	10.44	204	461481	38.73	nq	93
61) 4-Nitroaniline	10.48	138	340264	38.95	nq	# 80
62) Azobenzene	10.60	77	1304497	38.03	nq	95
64) 4,6-Dinitro-2-methylphenol	10.50	198	89722	39.07	nq	# 63
65) n-Nitrosodiphenylamine	10.57	169	989323	39.76	nq	99
66) 4-Bromophenyl-phenylether	10.93	248	299694	36.93	nq	# 87
67) Hexachlorobenzene	11.00	284	304951	37.44	nq	# 72
68) Atrazine	11.09	200	300511	39.29	nq	96
69) Pentachlorophenol	11.18	266	169740	39.79	nq	97
70) Phenanthrene	11.41	178	1440037	40.24	nq	99
71) Anthracene	11.46	178	1432667	37.38	nq	98
72) Carbazole	11.62	167	1389152	38.30	nq	98
73) Di-n-butylphthalate	11.95	149	1820990	39.00	nq	100
74) Fluoranthene	12.59	202	1207355	36.39	nq	98
76) Benzidine	12.73	184	713214	37.79	nq	99
77) Pyrene	12.82	202	1178882	37.18	nq	99
79) Butylbenzylphthalate	13.45	149	655611	39.17	nq	95
80) Benzo(a)anthracene	14.01	228	875200	38.87	nq	99
81) 3,3'-Dichlorobenzidine	13.97	252	591415	37.97	nq	97
82) Chrysene	14.04	228	875014	35.95	nq	99
83) Bis(2-ethylhexyl)phthalate	14.01	149	749473	38.24	nq	99
84) Di-n-octyl phthalate	14.61	149	1372791	38.86	nq	96
85) Indeno(1,2,3-cd)pyrene	16.84	276	1038425	41.49	nq	96
87) Benzo(b)fluoranthene	15.04	252	964794	35.43	nq	100
88) Benzo(k)fluoranthene	15.07	252	994757m	43.39	nq	
89) Benzo(a)pyrene	15.39	252	980707	39.07	nq	98
90) Dibenzo(a,h)anthracene	16.87	278	871334	39.77	nq	99
91) Benzo(g,h,i)perylene	17.27	276	837912	39.89	nq	97

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

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