

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020817\
 Data File : BF092872.D
 Acq On : 8 Feb 2017 23:58
 Operator : SJ/MA
 Sample : I1751-04MS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ROLL-OFF-COMPMS

Manual Integrations
 APPROVED

mohammad
 2/9/2017 12:31:31 PM

Quant Time: Feb 09 00:45:30 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 08 23:48:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	151514	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	584360	20.00	ng	-0.01
38) Acenaphthene-d10	9.94	164	337387	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	562758	20.00	ng	0.00
75) Chrysene-d12	14.07	240	495391	20.00	ng	0.00
86) Perylene-d12	15.51	264	366215	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.52	112	1112770	126.00	ng	0.02
7) Phenol-d6	6.54	99	1374252	120.94	ng	0.01
23) Nitrobenzene-d5	7.47	82	1035801	98.71	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	378344	155.05	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	1909171	102.52	ng	0.00
78) Terphenyl-d14	13.01	244	1826004	86.61	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.67	88	175428	36.76	ng	# 73
3) Pyridine	3.38	79	385873m	31.80	ng	
4) n-Nitrosodimethylamine	3.31	42	254936	44.74	ng	91
6) Aniline	6.57	93	126810	8.18	ng	# 81
8) 2-Chlorophenol	6.69	128	453725	46.57	ng	89
9) Benzaldehyde	6.44	77	113324	13.92	ng	89
10) Phenol	6.56	94	470668	35.40	ng	79
11) bis(2-Chloroethyl)ether	6.64	93	463041	43.64	ng	92
12) 1,3-Dichlorobenzene	6.84	146	518991	45.48	ng	98
13) 1,4-Dichlorobenzene	6.92	146	504258m	43.66	ng	
14) 1,2-Dichlorobenzene	7.07	146	495301	44.85	ng	97
15) Benzyl Alcohol	7.05	79	400338	47.59	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.18	45	844989	45.83	ng	96
17) 2-Methylphenol	7.16	107	395849	47.36	ng	96
18) Hexachloroethane	7.41	117	169278	42.93	ng	90
19) n-Nitroso-di-n-propylamine	7.32	70	348918	48.24	ng	# 94
20) 3+4-Methylphenols	7.32	107	478620	47.89	ng	# 86
22) Acetophenone	7.31	105	610629	45.77	ng	# 95
24) Nitrobenzene	7.48	77	469004	43.53	ng	99
25) Isophorone	7.72	82	934577	47.79	ng	96
26) 2-Nitrophenol	7.80	139	274631	53.62	ng	92
27) 2,4-Dimethylphenol	7.85	122	473501	52.64	ng	94
28) bis(2-Chloroethoxy)methane	7.94	93	634685	49.01	ng	99
29) 2,4-Dichlorophenol	8.04	162	417557	49.77	ng	88
30) 1,2,4-Trichlorobenzene	8.12	180	416976	46.12	ng	95
31) Naphthalene	8.20	128	1293453	43.66	ng	99
32) Benzoic acid	7.98	122	244022	47.90	ng	86
33) 4-Chloroaniline	8.26	127	58884	5.07	ng	96
34) Hexachlorobutadiene	8.32	225	220320	44.29	ng	100
35) Caprolactam	8.64	113	118920m	45.06	ng	
36) 4-Chloro-3-methylphenol	8.75	107	458084	52.37	ng	97
37) 2-Methylnaphthalene	8.90	142	977035	51.37	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.06	216	396279	41.84	ng	99
40) Hexachlorocyclopentadiene	9.05	237	326758	76.47	ng	96

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42) 2,4,6-Trichlorophenol	9.17	196	302488	48.61	nq	90
43) 2,4,5-Trichlorophenol	9.22	196	323220	47.40	nq	95
45) 1,1'-Biphenyl	9.36	154	1073492	43.94	nq	97
46) 2-Chloronaphthalene	9.39	162	942329	49.31	nq	96
47) 2-Nitroaniline	9.48	65	268669	41.81	nq	98
48) Acenaphthylene	9.80	152	1444218	43.74	nq	99
49) Dimethylphthalate	9.66	163	1086304	47.80	nq	99
50) 2,6-Dinitrotoluene	9.73	165	256998	52.37	nq #	70
51) Acenaphthene	9.97	154	925233	46.87	nq	96
52) 3-Nitroaniline	9.89	138	26092	4.65	nq #	79
53) 2,4-Dinitrophenol	10.01	184	249608	98.80	nq #	78
54) Dibenzofuran	10.14	168	1238478	46.91	nq	97
55) 4-Nitrophenol	10.08	139	391176	96.70	nq	83
56) 2,4-Dinitrotoluene	10.13	165	317708	48.62	nq	96
57) Fluorene	10.49	166	964312	47.48	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.27	232	250666	55.11	nq	96
59) Diethylphthalate	10.36	149	1043232	48.71	nq	99
60) 4-Chlorophenyl-phenylether	10.48	204	413619	42.39	nq	92
61) 4-Nitroaniline	10.51	138	185379	32.23	nq	93
62) Azobenzene	10.64	77	1114438	50.37	nq	97
64) 4,6-Dinitro-2-methylphenol	10.54	198	195910	56.68	nq	93
65) n-Nitrosodiphenylamine	10.60	169	829524	46.14	nq	99
66) 4-Bromophenyl-phenylether	10.97	248	273527	46.52	nq	96
67) Hexachlorobenzene	11.04	284	285471	49.13	nq	96
68) Atrazine	11.13	200	218942	37.95	nq	97
69) Pentachlorophenol	11.23	266	295450	96.84	nq	97
70) Phenanthrene	11.45	178	1401625	43.47	nq	98
71) Anthracene	11.50	178	1631045	50.38	nq	98
72) Carbazole	11.65	167	1123585	36.30	nq	99
73) Di-n-butylphthalate	11.99	149	1791525	53.52	nq #	97
74) Fluoranthene	12.64	202	1511344	45.44	nq	97
77) Pyrene	12.87	202	1528363	41.23	nq	99
79) Butylbenzylphthalate	13.48	149	785972	52.21	nq	87
80) Benzo(a)anthracene	14.05	228	1355099	44.72	nq	99
81) 3,3'-Dichlorobenzidine	14.01	252	58481	5.41	nq	96
82) Chrysene	14.09	228	1073853	37.85	nq	99
83) Bis(2-ethylhexyl)phthalate	14.04	149	1048847	50.94	nq #	95
84) Di-n-octyl phthalate	14.65	149	1737422	51.82	nq	100
85) Indeno(1,2,3-cd)pyrene	16.95	276	1150010	52.34	nq	95
87) Benzo(b)fluoranthene	15.09	252	1262337	52.37	nq	99
88) Benzo(k)fluoranthene	15.13	252	1182389m	56.81	nq	
89) Benzo(a)pyrene	15.45	252	1143689	54.85	nq	97
90) Dibenzo(a,h)anthracene	16.97	278	962030	57.09	nq	98
91) Benzo(g,h,i)perylene	17.38	276	974060	57.22	nq #	91

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

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