

Data Path : Z:\HPCHEM1\BNA F\DATA\BF082416\
 Data File : BF089878.D
 Acq On : 24 Aug 2016 16:55
 Operator : UM/SJ
 Sample : PB93061BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB93061BS

Manual Integrations
 APPROVED

Quant Time: Aug 25 13:50:41 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 19 14:02:57 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	409033	20.00	ng	-0.02
21) Naphthalene-d8	7.95	136	1645721	20.00	ng	-0.02
38) Acenaphthene-d10	9.71	164	788148	20.00	ng	0.00
63) Phenanthrene-d10	11.19	188	1420523	20.00	ng	0.00
75) Chrysene-d12	13.84	240	1043361	20.00	ng	-0.03
86) Perylene-d12	15.31	264	910283	20.00	ng	-0.06

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System Monitoring Compounds

5) 2-Fluorophenol	5.24	112	2856219	119.75	ng	-0.01
7) Phenol-d6	6.31	99	3199601	109.28	ng	-0.01
23) Nitrobenzene-d5	7.23	82	1893144	71.47	ng	-0.02
41) 2,4,6-Tribromophenol	10.49	330	801700	107.03	ng	-0.01
44) 2-Fluorobiphenyl	9.04	172	3678568	82.03	ng	-0.01
78) Terphenyl-d14	12.78	244	3132809	67.93	ng	-0.01

8/25/2016 6:49:28 PM

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.20	88	386917	32.78	ng	95
3) Pyridine	2.83	79	1165020	37.63	ng	91
4) n-Nitrosodimethylamine	2.79	42	437755	37.98	ng	# 85
6) Aniline	6.33	93	1143292	29.11	ng	99
8) 2-Chlorophenol	6.44	128	1043866	36.60	ng	99
9) Benzaldehyde	6.20	77	240264	12.60	ng	87
10) Phenol	6.32	94	1197494	34.90	ng	99
11) bis(2-Chloroethyl)ether	6.41	93	1069142	40.18	ng	99
12) 1,3-Dichlorobenzene	6.60	146	1121541	37.61	ng	98
13) 1,4-Dichlorobenzene	6.68	146	1185151	38.66	ng	98
14) 1,2-Dichlorobenzene	6.84	146	1056110	36.91	ng	98
15) Benzyl Alcohol	6.81	79	822730	42.37	ng	100
16) 2,2'-oxybis(1-Chloropropan	6.96	45	1139563	33.46	ng	93
17) 2-Methylphenol	6.94	107	939662	42.06	ng	95
18) Hexachloroethane	7.19	117	419295	38.35	ng	90
19) n-Nitroso-di-n-propylamine	7.10	70	761488	40.67	ng	# 88
20) 3+4-Methylphenols	7.08	107	1122332	40.97	ng	94
22) Acetophenone	7.08	105	1314575	34.79	ng	# 90
24) Nitrobenzene	7.26	77	1198486	40.32	ng	91
25) Isophorone	7.50	82	1974779	36.96	ng	96
26) 2-Nitrophenol	7.58	139	597699	40.31	ng	98
27) 2,4-Dimethylphenol	7.62	122	1086572	40.66	ng	99
28) bis(2-Chloroethoxy)methane	7.71	93	1186062	35.88	ng	98
29) 2,4-Dichlorophenol	7.82	162	949376	42.34	ng	99
30) 1,2,4-Trichlorobenzene	7.90	180	906908	36.56	ng	99
31) Naphthalene	7.98	128	2965251	38.55	ng	99
32) Benzoic acid	7.74	122	532098	25.55	ng	97
33) 4-Chloroaniline	8.02	127	564627	16.94	ng	# 86
34) Hexachlorobutadiene	8.10	225	480402	36.98	ng	99
35) Caprolactam	8.40	113	280872	41.13	ng	# 80
36) 4-Chloro-3-methylphenol	8.51	107	928512	38.23	ng	90
37) 2-Methylnaphthalene	8.67	142	2075832	41.72	ng	100
39) 1,2,4,5-Tetrachlorobenzene	8.83	216	670854	26.29	ng	99
40) Hexachlorocyclopentadiene	8.83	237	850548	95.99	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

42) 2,4,6-Trichlorophenol	8.95	196	631651	39.34	nq	100
43) 2,4,5-Trichlorophenol	8.99	196	632562	39.61	nq	# 93
45) 1,1'-Biphenyl	9.13	154	2068188	33.91	nq	99
46) 2-Chloronaphthalene	9.15	162	1786734	37.06	nq	98
47) 2-Nitroaniline	9.26	65	613622	44.62	nq	# 64
48) Acenaphthylene	9.56	152	2850808	39.35	nq	99
49) Dimethylphthalate	9.45	163	2172944	39.16	nq	# 98
50) 2,6-Dinitrotoluene	9.50	165	461198	36.00	nq	87
51) Acenaphthene	9.74	154	1991746	41.54	nq	99
52) 3-Nitroaniline	9.66	138	343835	24.22	nq	97
53) 2,4-Dinitrophenol	9.77	184	456106	65.66	nq	# 78
54) Dibenzofuran	9.91	168	2604258	43.24	nq	98
55) 4-Nitrophenol	9.83	139	773000	66.72	nq	90
56) 2,4-Dinitrotoluene	9.90	165	602994	36.27	nq	# 64
57) Fluorene	10.25	166	1724218	35.93	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.03	232	550372	47.35	nq	97
59) Diethylphthalate	10.15	149	2228472	39.53	nq	98
60) 4-Chlorophenyl-phenylether	10.25	204	777671	35.52	nq	95
61) 4-Nitroaniline	10.27	138	480472	36.10	nq	95
62) Azobenzene	10.41	77	2141606	40.41	nq	95
64) 4,6-Dinitro-2-methylphenol	10.31	198	287666	35.24	nq	# 58
65) n-Nitrosodiphenylamine	10.38	169	1857356	41.58	nq	100
66) 4-Bromophenyl-phenylether	10.74	248	580043	38.95	nq	# 89
67) Hexachlorobenzene	10.80	284	619348	36.45	nq	# 87
68) Atrazine	10.90	200	486543	35.57	nq	97
69) Pentachlorophenol	10.99	266	662465	68.79	nq	97
70) Phenanthrene	11.21	178	2966461	41.35	nq	97
71) Anthracene	11.26	178	2699943	37.02	nq	99
72) Carbazole	11.42	167	2723684	41.64	nq	97
73) Di-n-butylphthalate	11.77	149	3346838	40.89	nq	# 98
74) Fluoranthene	12.39	202	2671296	39.14	nq	94
76) Benzidine	12.51	184	1151668	34.11	nq	97
77) Pyrene	12.62	202	2846865	33.51	nq	97
79) Butylbenzylphthalate	13.27	149	1575622	41.06	nq	93
80) Benzo(a)anthracene	13.83	228	2476214	41.44	nq	99
81) 3,3'-Dichlorobenzidine	13.81	252	644010	32.88	nq	# 96
82) Chrysene	13.87	228	2201057	42.97	nq	99
83) Bis(2-ethylhexyl)phthalate	13.86	149	2059673	38.96	nq	97
84) Di-n-octyl phthalate	14.53	149	3473434	49.45	nq	# 93
85) Indeno(1,2,3-cd)pyrene	16.61	276	2032387	31.10	nq	99
87) Benzo(b)fluoranthene	14.94	252	2359067	47.11	nq	# 95
88) Benzo(k)fluoranthene	14.96	252	1724474m	37.95	nq	
89) Benzo(a)pyrene	15.26	252	1904059	40.17	nq	99
90) Dibenzo(a,h)anthracene	16.63	278	1684446	32.74	nq	97
91) Benzo(g,h,i)perylene	16.98	276	1715358	31.66	nq	100

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

