

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030817\
 Data File : BF093432.D
 Acq On : 8 Mar 2017 17:01
 Operator : SJ/MA
 Sample : PB97430BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB97430BS

Manual Integrations
 APPROVED

mohammad
 3/10/2017 8:11:39 AM

Quant Time: Mar 08 17:41:33 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 07 15:55:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	192739m	20.00	ng	-0.01
21) Naphthalene-d8	8.05	136	766439	20.00	ng	-0.01
38) Acenaphthene-d10	9.81	164	382551	20.00	ng	0.00
63) Phenanthrene-d10	11.29	188	706812	20.00	ng	0.01
75) Chrysene-d12	13.93	240	622673	20.00	ng	0.01
86) Perylene-d12	15.34	264	496995	20.00	ng	0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	1272029	109.84	ng	0.01
7) Phenol-d6	6.42	99	1747585	119.67	ng	0.01
23) Nitrobenzene-d5	7.34	82	1058821	76.65	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	320296	128.51	ng	0.00
44) 2-Fluorobiphenyl	9.13	172	1814525	88.23	ng	0.00
78) Terphenyl-d14	12.87	244	1502048	62.72	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.33	88	182837	28.66	ng	# 82
3) Pyridine	3.03	79	530430	32.62	ng	87
4) n-Nitrosodimethylamine	2.98	42	273948	35.27	ng	84
6) Aniline	6.44	93	259777m	12.96	ng	
8) 2-Chlorophenol	6.55	128	469867	36.21	ng	95
9) Benzaldehyde	6.32	77	79940	6.35	ng	98
10) Phenol	6.44	94	607601m	33.95	ng	
11) bis(2-Chloroethyl)ether	6.50	93	503061	34.78	ng	96
12) 1,3-Dichlorobenzene	6.70	146	504168m	33.95	ng	
13) 1,4-Dichlorobenzene	6.78	146	523742m	34.44	ng	
14) 1,2-Dichlorobenzene	6.94	146	452693	33.62	ng	94
15) Benzyl Alcohol	6.93	79	380859	36.59	ng	93
16) 2,2'-oxybis(1-Chloropropan	7.05	45	823404	34.27	ng	85
17) 2-Methylphenol	7.04	107	394983	35.99	ng	94
18) Hexachloroethane	7.27	117	177613	34.64	ng	99
19) n-Nitroso-di-n-propylamine	7.19	70	378486	37.70	ng	97
20) 3+4-Methylphenols	7.20	107	546676	38.41	ng	# 73
22) Acetophenone	7.18	105	656112	35.57	ng	# 97
24) Nitrobenzene	7.36	77	546624	35.21	ng	# 89
25) Isophorone	7.60	82	990088	35.54	ng	99
26) 2-Nitrophenol	7.67	139	257209	36.31	ng	# 87
27) 2,4-Dimethylphenol	7.73	122	428818	37.21	ng	92
28) bis(2-Chloroethoxy)methane	7.81	93	627255	35.67	ng	99
29) 2,4-Dichlorophenol	7.92	162	423329	39.05	ng	88
30) 1,2,4-Trichlorobenzene	7.99	180	404417	35.08	ng	# 95
31) Naphthalene	8.07	128	1292831	33.66	ng	99
32) Benzoic acid	7.89	122	293916	49.09	ng	91
33) 4-Chloroaniline	8.14	127	170974	10.97	ng	98
34) Hexachlorobutadiene	8.18	225	211818	35.67	ng	99
35) Caprolactam	8.52	113	149603m	40.41	ng	
36) 4-Chloro-3-methylphenol	8.63	107	439959	38.02	ng	92
37) 2-Methylnaphthalene	8.77	142	923745	37.80	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.93	216	380536	36.43	ng	99
40) Hexachlorocyclopentadiene	8.92	237	206719	78.66	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.05	196	271622	41.62	nq	93
43) 2,4,5-Trichlorophenol	9.11	196	261602	37.36	nq	# 86
45) 1,1'-Biphenyl	9.22	154	1026691	36.84	nq	96
46) 2-Chloronaphthalene	9.26	162	871162	40.84	nq	95
47) 2-Nitroaniline	9.36	65	309478	41.04	nq	95
48) Acenaphthylene	9.67	152	1332146	37.87	nq	98
49) Dimethylphthalate	9.53	163	1006773	39.70	nq	99
50) 2,6-Dinitrotoluene	9.60	165	205511	36.93	nq	# 52
51) Acenaphthene	9.84	154	818164	39.33	nq	98
52) 3-Nitroaniline	9.77	138	98549	14.74	nq	91
53) 2,4-Dinitrophenol	9.90	184	217410	85.79	nq	# 67
54) Dibenzofuran	10.01	168	1151724	44.23	nq	94
55) 4-Nitrophenol	9.97	139	280151	86.27	nq	94
56) 2,4-Dinitrotoluene	10.01	165	266311	38.95	nq	# 65
57) Fluorene	10.36	166	909863	41.24	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.14	232	218135	44.14	nq	97
59) Diethylphthalate	10.23	149	931944	38.45	nq	100
60) 4-Chlorophenyl-phenylether	10.34	204	416540	42.12	nq	88
61) 4-Nitroaniline	10.40	138	249541	36.49	nq	79
62) Azobenzene	10.50	77	1218109	44.23	nq	96
64) 4,6-Dinitro-2-methylphenol	10.44	198	162391	35.46	nq	89
65) n-Nitrosodiphenylamine	10.47	169	820854	34.78	nq	99
66) 4-Bromophenyl-phenylether	10.84	248	259491	34.72	nq	# 83
67) Hexachlorobenzene	10.90	284	250668	35.12	nq	97
68) Atrazine	11.00	200	227503	32.93	nq	99
69) Pentachlorophenol	11.11	266	213708	86.73	nq	98
70) Phenanthrene	11.32	178	1394393	34.56	nq	99
71) Anthracene	11.36	178	1414175	34.65	nq	99
72) Carbazole	11.53	167	1310558	34.18	nq	# 96
73) Di-n-butylphthalate	11.85	149	1731428	39.04	nq	# 96
74) Fluoranthene	12.50	202	1450823	37.23	nq	95
76) Benzidine	12.63	184	305193	14.90	nq	96
77) Pyrene	12.73	202	1462042	32.28	nq	99
79) Butylbenzylphthalate	13.35	149	732926	38.45	nq	# 77
80) Benzo(a)anthracene	13.92	228	1223794	33.28	nq	99
81) 3,3'-Dichlorobenzidine	13.89	252	285596	22.18	nq	# 96
82) Chrysene	13.96	228	1031834	30.33	nq	100
83) Bis(2-ethylhexyl)phthalate	13.90	149	956975	36.02	nq	100
84) Di-n-octyl phthalate	14.52	149	1692493	38.22	nq	98
85) Indeno(1,2,3-cd)pyrene	16.72	276	893976	31.39	nq	# 93
87) Benzo(b)fluoranthene	14.95	252	1249299m	41.27	nq	
88) Benzo(k)fluoranthene	14.97	252	857769m	29.39	nq	
89) Benzo(a)pyrene	15.29	252	997842	36.07	nq	96
90) Dibenzo(a,h)anthracene	16.72	278	756682	33.06	nq	95
91) Benzo(g,h,i)perylene	17.13	276	771283	32.84	nq	# 93

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

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