

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012916\
 Data File : BF084648.D
 Acq On : 29 Jan 2016 21:19
 Operator : SJ/IZ
 Sample : PB88084BS
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB88084BS

Manual Integrations
 APPROVED

mohammad
 2/1/2016 2:24:30 PM

Quant Time: Jan 30 00:15:17 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 29 12:54:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	172185	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	687774	20.00	ng	-0.01
38) Acenaphthene-d10	9.77	164	329859	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	634360	20.00	ng	0.00
75) Chrysene-d12	13.88	240	514077	20.00	ng	0.01
86) Perylene-d12	15.25	264	518561	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.33	112	1354744	118.77	ng	0.02
7) Phenol-d6	6.37	99	1562485	114.29	ng	0.00
23) Nitrobenzene-d5	7.30	82	1075736	87.41	ng	0.00
41) 2,4,6-Tribromophenol	10.56	330	457701	132.39	ng	0.00
44) 2-Fluorobiphenyl	9.09	172	1773592	79.92	ng	0.00
78) Terphenyl-d14	12.83	244	1802653	79.20	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.30	88	163672	30.77	ng	# 77
3) Pyridine	2.98	79	426210	30.08	ng	# 77
4) n-Nitrosodimethylamine	2.93	42	254445	36.72	ng	81
6) Aniline	6.38	93	330566	18.84	ng	# 1
8) 2-Chlorophenol	6.51	128	499074	39.44	ng	91
9) Benzaldehyde	6.27	77	216276	27.32	ng	90
10) Phenol	6.38	94	509380	33.43	ng	81
11) bis(2-Chloroethyl)ether	6.46	93	430898	36.36	ng	# 80
12) 1,3-Dichlorobenzene	6.66	146	475721m	35.09	ng	
13) 1,4-Dichlorobenzene	6.74	146	491191	36.92	ng	97
14) 1,2-Dichlorobenzene	6.90	146	478752	39.08	ng	96
15) Benzyl Alcohol	6.88	79	391048	42.02	ng	92
16) 2,2'-oxybis(1-Chloropropan	7.01	45	730580	36.02	ng	86
17) 2-Methylphenol	7.00	107	405507	41.34	ng	94
18) Hexachloroethane	7.24	117	196456	37.75	ng	# 82
19) n-Nitroso-di-n-propylamine	7.15	70	343475	41.13	ng	# 77
20) 3+4-Methylphenols	7.15	107	488370	41.03	ng	# 78
22) Acetophenone	7.14	105	673387	37.49	ng	98
24) Nitrobenzene	7.31	77	484763	37.06	ng	96
25) Isophorone	7.55	82	932979	40.46	ng	97
26) 2-Nitrophenol	7.63	139	266758	42.88	ng	# 87
27) 2,4-Dimethylphenol	7.68	122	446813	40.99	ng	94
28) bis(2-Chloroethoxy)methane	7.77	93	553556	40.00	ng	100
29) 2,4-Dichlorophenol	7.88	162	402592	42.23	ng	94
30) 1,2,4-Trichlorobenzene	7.95	180	412051	37.79	ng	96
31) Naphthalene	8.03	128	1293895	37.27	ng	99
32) Benzoic acid	7.81	122	299700	35.82	ng	97
33) 4-Chloroaniline	8.09	127	134135	9.50	ng	96
34) Hexachlorobutadiene	8.16	225	253096	39.28	ng	99
35) Caprolactam	8.46	113	138181m	50.51	ng	
36) 4-Chloro-3-methylphenol	8.58	107	470115	44.79	ng	93
37) 2-Methylnaphthalene	8.73	142	984410	46.19	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.89	216	371640	34.74	ng	98
40) Hexachlorocyclopentadiene	8.88	237	423166	87.00	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.01	196	277575	41.77	nq	92
43) 2,4,5-Trichlorophenol	9.05	196	275566	40.16	nq	# 80
45) 1,1'-Biphenyl	9.20	154	1126631	37.27	nq	95
46) 2-Chloronaphthalene	9.21	162	836879	37.94	nq	96
47) 2-Nitroaniline	9.31	65	267329	40.36	nq	98
48) Acenaphthylene	9.63	152	1363772	40.81	nq	98
49) Dimethylphthalate	9.50	163	1053993	42.57	nq	# 91
50) 2,6-Dinitrotoluene	9.56	165	242988	44.61	nq	98
51) Acenaphthene	9.80	154	896893	40.08	nq	98
52) 3-Nitroaniline	9.72	138	87158	15.31	nq	# 88
53) 2,4-Dinitrophenol	9.84	184	270560	100.66	nq	91
54) Dibenzofuran	9.97	168	1254210	46.96	nq	97
55) 4-Nitrophenol	9.90	139	370362	88.58	nq	# 83
56) 2,4-Dinitrotoluene	9.96	165	322144	45.52	nq	# 93
57) Fluorene	10.32	166	948476	41.90	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.09	232	247796	47.77	nq	89
59) Diethylphthalate	10.20	149	1056614	43.60	nq	97
60) 4-Chlorophenyl-phenylether	10.30	204	394515	38.82	nq	92
61) 4-Nitroaniline	10.34	138	228671	42.44	nq	93
62) Azobenzene	10.46	77	1055639	45.51	nq	90
64) 4,6-Dinitro-2-methylphenol	10.37	198	177194	42.77	nq	70
65) n-Nitrosodiphenylamine	10.43	169	877051	42.64	nq	99
66) 4-Bromophenyl-phenylether	10.80	248	304753	42.66	nq	96
67) Hexachlorobenzene	10.85	284	293427	37.76	nq	# 83
68) Atrazine	10.96	200	255905	41.52	nq	98
69) Pentachlorophenol	11.06	266	318283	84.77	nq	100
70) Phenanthrene	11.26	178	1345974	38.37	nq	97
71) Anthracene	11.32	178	1544625	43.31	nq	98
72) Carbazole	11.48	167	1422273	46.27	nq	98
73) Di-n-butylphthalate	11.81	149	1456685	38.97	nq	# 97
74) Fluoranthene	12.45	202	1611709	46.47	nq	95
76) Benzidine	12.58	184	626940	32.07	nq	98
77) Pyrene	12.68	202	1653404	42.64	nq	98
79) Butylbenzylphthalate	13.31	149	751227	45.35	nq	# 94
80) Benzo(a)anthracene	13.86	228	1354013	42.25	nq	99
81) 3,3'-Dichlorobenzidine	13.84	252	195015	17.30	nq	100
82) Chrysene	13.91	228	1384891	43.84	nq	99
83) Bis(2-ethylhexyl)phthalate	13.87	149	906889	42.21	nq	98
84) Di-n-octyl phthalate	14.48	149	1624294	47.84	nq	# 90
85) Indeno(1,2,3-cd)pyrene	16.57	276	1134553	40.63	nq	99
87) Benzo(b)fluoranthene	14.88	252	1322203m	39.38	nq	
88) Benzo(k)fluoranthene	14.90	252	1175430m	42.38	nq	
89) Benzo(a)pyrene	15.21	252	1247048	42.93	nq	99
90) Dibenzo(a,h)anthracene	16.59	278	956010	36.52	nq	# 95
91) Benzo(g,h,i)perylene	16.97	276	907317	34.33	nq	97

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

