

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112916\
 Data File : BF091320.D
 Acq On : 29 Nov 2016 11:53
 Operator : UM/SJ
 Sample : SSTDICC050
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

sohil
 11/30/2016 10:32:11 AM

Quant Time: Nov 29 12:30:01 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 29 12:14:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	170909	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	646243	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	381858	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	634683	20.00	ng	0.00
75) Chrysene-d12	14.03	240	424295	20.00	ng	0.01
86) Perylene-d12	15.44	264	371362	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	1060268	99.67	ng	0.00
7) Phenol-d6	6.49	99	1240369	90.76	ng	0.00
23) Nitrobenzene-d5	7.43	82	1122565	100.00	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	371478	96.47	ng	0.01
44) 2-Fluorobiphenyl	9.24	172	2129089	91.30	ng	0.00
78) Terphenyl-d14	12.97	244	1886539	101.55	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.47	88	236518	47.14	ng	88
3) Pyridine	3.19	79	668447	49.57	ng	# 85
4) n-Nitrosodimethylamine	3.14	42	361703	48.99	ng	98
6) Aniline	6.53	93	836002	47.85	ng	# 71
8) 2-Chlorophenol	6.65	128	571271	49.30	ng	93
9) Benzaldehyde	6.41	77	398633	53.89	ng	83
10) Phenol	6.50	94	688794	45.17	ng	97
11) bis(2-Chloroethyl)ether	6.61	93	573007	51.12	ng	95
12) 1,3-Dichlorobenzene	6.81	146	610293m	48.31	ng	
13) 1,4-Dichlorobenzene	6.88	146	583343	45.11	ng	96
14) 1,2-Dichlorobenzene	7.04	146	613048	49.97	ng	96
15) Benzyl Alcohol	7.01	79	474708	44.73	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.14	45	1150858	46.83	ng	71
17) 2-Methylphenol	7.12	107	432813	47.21	ng	# 76
18) Hexachloroethane	7.38	117	224512	48.77	ng	96
19) n-Nitroso-di-n-propylamine	7.29	70	412958	48.17	ng	# 84
20) 3+4-Methylphenols	7.28	107	559676	49.73	ng	# 41
22) Acetophenone	7.28	105	754106	51.71	ng	# 85
24) Nitrobenzene	7.45	77	665893	56.06	ng	# 88
25) Isophorone	7.69	82	1004018	47.88	ng	99
26) 2-Nitrophenol	7.77	139	286499	60.14	ng	# 70
27) 2,4-Dimethylphenol	7.81	122	518248	51.84	ng	96
28) bis(2-Chloroethoxy)methane	7.91	93	618024	49.50	ng	# 96
29) 2,4-Dichlorophenol	8.01	162	445472	50.42	ng	90
30) 1,2,4-Trichlorobenzene	8.09	180	484125	47.57	ng	99
31) Naphthalene	8.17	128	1449664	46.63	ng	100
32) Benzoic acid	7.93	122	404067	51.32	ng	93
33) 4-Chloroaniline	8.22	127	643629	47.59	ng	# 93
34) Hexachlorobutadiene	8.30	225	309380	51.42	ng	98
35) Caprolactam	8.60	113	141492	51.95	ng	98
36) 4-Chloro-3-methylphenol	8.70	107	464254	48.89	ng	100
37) 2-Methylnaphthalene	8.87	142	1080167	50.63	ng	95
39) 1,2,4,5-Tetrachlorobenzene	9.03	216	445973	43.10	ng	98
40) Hexachlorocyclopentadiene	9.02	237	237942	51.37	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.14	196	349256	48.94	ng	97
43) 2,4,5-Trichlorophenol	9.18	196	337058	49.16	ng #	90
45) 1,1'-Biphenyl	9.33	154	1213805	44.09	ng	98
46) 2-Chloronaphthalene	9.35	162	944231	46.82	ng	98
47) 2-Nitroaniline	9.44	65	344329	50.32	ng #	74
48) Acenaphthylene	9.77	152	1629852	50.43	ng	99
49) Dimethylphthalate	9.64	163	1077524	45.19	ng	98
50) 2,6-Dinitrotoluene	9.69	165	281425	54.04	ng #	82
51) Acenaphthene	9.94	154	1065693	48.45	ng	97
52) 3-Nitroaniline	9.86	138	314643	56.19	ng #	77
53) 2,4-Dinitrophenol	9.96	184	117123	62.00	ng #	65
54) Dibenzofuran	10.12	168	1424537	46.04	ng	93
55) 4-Nitrophenol	10.01	139	251242	62.21	ng	90
56) 2,4-Dinitrotoluene	10.09	165	361423	54.84	ng #	79
57) Fluorene	10.46	166	1123415	47.45	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.23	232	308962	51.24	ng	95
59) Diethylphthalate	10.33	149	1084902	45.22	ng	96
60) 4-Chlorophenyl-phenylether	10.45	204	461048	42.27	ng #	85
61) 4-Nitroaniline	10.47	138	259002	49.58	ng	89
62) Azobenzene	10.61	77	1161024	47.71	ng	89
64) 4,6-Dinitro-2-methylphenol	10.50	198	196965	68.43	ng #	34
65) n-Nitrosodiphenylamine	10.56	169	887883	46.08	ng	99
66) 4-Bromophenyl-phenylether	10.94	248	337748	48.69	ng #	86
67) Hexachlorobenzene	11.01	284	356527	46.88	ng #	88
68) Atrazine	11.10	200	311602	61.26	ng	92
69) Pentachlorophenol	11.19	266	225080	50.03	ng	96
70) Phenanthrene	11.42	178	1593484	48.40	ng	99
71) Anthracene	11.46	178	1547550	45.97	ng	99
72) Carbazole	11.61	167	1351934	45.12	ng	99
73) Di-n-butylphthalate	11.96	149	1647741	47.49	ng	100
74) Fluoranthene	12.60	202	1560082	46.58	ng	98
76) Benzidine	12.72	184	678106	82.46	ng	98
77) Pyrene	12.82	202	1602982	51.13	ng	99
79) Butylbenzylphthalate	13.45	149	657506	52.41	ng #	82
80) Benzo(a)anthracene	14.01	228	1247680	51.81	ng	98
81) 3,3'-Dichlorobenzidine	13.98	252	449884	56.65	ng #	98
82) Chrysene	14.05	228	1289517	56.66	ng	99
83) Bis(2-ethylhexyl)phthalate	14.01	149	819550	52.47	ng #	94
84) Di-n-octyl phthalate	14.62	149	1182765	48.11	ng	97
85) Indeno(1,2,3-cd)pyrene	16.86	276	1125115	57.41	ng	96
87) Benzo(b)fluoranthene	15.04	252	1198064	51.14	ng #	95
88) Benzo(k)fluoranthene	15.07	252	888243m	46.42	ng	
89) Benzo(a)pyrene	15.39	252	1012904	50.82	ng	97
90) Dibenzo(a,h)anthracene	16.87	278	947612	51.50	ng	96
91) Benzo(g,h,i)perylene	17.28	276	992724	54.39	ng	98

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

