

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062915\
 Data File : BF080220.D
 Acq On : 29 Jun 2015 23:22
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

TEJASKUMAR
 6/30/2015 2:18:57 PM

Quant Time: Jun 30 11:13:26 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 23 00:53:56 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.29	152	47146	20.00	ng	-0.03
21) Naphthalene-d8	8.58	136	191180	20.00	ng	-0.03
38) Acenaphthene-d10	10.36	164	100049	20.00	ng	-0.03
63) Phenanthrene-d10	11.86	188	208506	20.00	ng	-0.05
75) Chrysene-d12	14.80	240	194500	20.00	ng	-0.19
86) Perylene-d12	16.67	264	196808	20.00	ng	-0.26

System Monitoring Compounds

5) 2-Fluorophenol	5.90	112	212653	79.84	ng	-0.03
7) Phenol-d6	6.90	99	291106	82.14	ng	-0.02
23) Nitrobenzene-d5	7.85	82	279002	84.96	ng	-0.02
41) 2,4,6-Tribromophenol	11.16	330	80853	82.64	ng	-0.02
44) 2-Fluorobiphenyl	9.66	172	511038	72.84	ng	-0.03
78) Terphenyl-d14	13.54	244	681651	81.85	ng	-0.13

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.12	88	44878m	35.45	ng	
3) Pyridine	3.96	79	115908	34.43	ng	# 82
4) n-Nitrosodimethylamine	3.86	42	64436	40.27	ng	95
6) Aniline	6.95	93	209956	43.06	ng	92
8) 2-Chlorophenol	7.08	128	127112	41.29	ng	89
9) Benzaldehyde	6.84	77	73259	35.81	ng	# 70
10) Phenol	6.92	94	168124	43.47	ng	83
11) bis(2-Chloroethyl)ether	7.02	93	111921	36.47	ng	85
12) 1,3-Dichlorobenzene	7.24	146	147361	39.85	ng	# 92
13) 1,4-Dichlorobenzene	7.32	146	151031	42.95	ng	94
14) 1,2-Dichlorobenzene	7.46	146	132771m	38.80	ng	
15) Benzyl Alcohol	7.42	79	111362	38.43	ng	# 74
16) 2,2'-oxybis(1-Chloropropan	7.57	45	199204	43.74	ng	84
17) 2-Methylphenol	7.53	107	107194	40.77	ng	# 89
18) Hexachloroethane	7.82	117	50010	42.71	ng	# 80
19) n-Nitroso-di-n-propylamine	7.69	70	105243	42.77	ng	# 77
20) 3+4-Methylphenols	7.68	107	145442	42.86	ng	90
22) Acetophenone	7.69	105	172264	38.67	ng	# 80
24) Nitrobenzene	7.88	77	142782	42.12	ng	# 77
25) Isophorone	8.10	82	247793	37.49	ng	# 83
26) 2-Nitrophenol	8.20	139	69730	49.15	ng	# 78
27) 2,4-Dimethylphenol	8.22	122	125201	42.32	ng	# 83
28) bis(2-Chloroethoxy)methane	8.31	93	146997	37.86	ng	# 91
29) 2,4-Dichlorophenol	8.44	162	113557	44.82	ng	100
30) 1,2,4-Trichlorobenzene	8.53	180	133099	46.26	ng	98
31) Naphthalene	8.61	128	394070m	39.42	ng	
32) Benzoic acid	8.29	122	55869	31.23	ng	# 64
33) 4-Chloroaniline	8.64	127	151360	35.38	ng	# 82
34) Hexachlorobutadiene	8.73	225	70382	39.73	ng	100
35) Caprolactam	9.00	113	36642	44.56	ng	94
36) 4-Chloro-3-methylphenol	9.12	107	114583	40.10	ng	92
37) 2-Methylnaphthalene	9.30	142	279185	43.18	ng	# 92
39) 1,2,4,5-Tetrachlorobenzene	9.48	216	126413	43.42	ng	98
40) Hexachlorocyclopentadiene	9.46	237	44317	33.52	ng	98

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42) 2,4,6-Trichlorophenol	9.58	196	88704	44.78	ng	99
43) 2,4,5-Trichlorophenol	9.61	196	87157	38.19	ng #	87
45) 1,1'-Biphenyl	9.76	154	304882	37.45	ng	91
46) 2-Chloronaphthalene	9.80	162	263194	39.85	ng	93
47) 2-Nitroaniline	9.88	65	72093	39.76	ng #	57
48) Acenaphthylene	10.22	152	458352	41.58	ng	97
49) Dimethylphthalate	10.06	163	311730	41.44	ng #	97
50) 2,6-Dinitrotoluene	10.12	165	60086	39.85	ng #	38
51) Acenaphthene	10.39	154	264451	39.21	ng	89
52) 3-Nitroaniline	10.30	138	72272	41.54	ng	97
53) 2,4-Dinitrophenol	10.40	184	24323	42.57	ng #	1
54) Dibenzofuran	10.56	168	363254	37.96	ng #	89
55) 4-Nitrophenol	10.44	139	52297	37.61	ng #	37
56) 2,4-Dinitrotoluene	10.53	165	88354	46.17	ng #	79
57) Fluorene	10.92	166	299065	39.39	ng	91
58) 2,3,4,6-Tetrachlorophenol	10.68	232	73423	38.11	ng	98
59) Diethylphthalate	10.76	149	284721	37.59	ng	97
60) 4-Chlorophenyl-phenylether	10.89	204	142105	38.94	ng #	88
61) 4-Nitroaniline	10.90	138	69491	38.67	ng #	41
62) Azobenzene	11.05	77	296687	38.48	ng	88
64) 4,6-Dinitro-2-methylphenol	10.94	198	35704	37.29	ng #	74
65) n-Nitrosodiphenylamine	11.01	169	256835	37.83	ng	93
66) 4-Bromophenyl-phenylether	11.40	248	86246	38.90	ng #	84
67) Hexachlorobenzene	11.48	284	93904	38.11	ng #	77
68) Atrazine	11.53	200	86969	40.54	ng	83
69) Pentachlorophenol	11.66	266	47631	34.11	ng	97
70) Phenanthrene	11.89	178	474863	37.62	ng	97
71) Anthracene	11.94	178	460850	37.92	ng	99
72) Carbazole	12.09	167	441688	38.06	ng	96
73) Di-n-butylphthalate	12.42	149	446278	33.29	ng	98
74) Fluoranthene	13.14	202	495456	36.85	ng	89
76) Benzidine	13.26	184	240974	44.32	ng	98
77) Pyrene	13.40	202	495680	41.39	ng	97
79) Butylbenzylphthalate	14.08	149	188443	39.19	ng	89
80) Benzo(a)anthracene	14.79	228	450507	38.02	ng	95
81) 3,3'-Dichlorobenzidine	14.73	252	152666	37.36	ng #	96
82) Chrysene	14.84	228	426580	39.04	ng	94
83) Bis(2-ethylhexyl)phthalate	14.77	149	270633	35.61	ng #	93
84) Di-n-octyl phthalate	15.57	149	438756	33.69	ng	95
85) Indeno(1,2,3-cd)pyrene	18.51	276	558571	40.54	ng #	88
87) Benzo(b)fluoranthene	16.13	252	469607	39.41	ng #	89
88) Benzo(k)fluoranthene	16.16	252	446059	36.76	ng #	91
89) Benzo(a)pyrene	16.59	252	445636	39.38	ng #	88
90) Dibenzo(a,h)anthracene	18.53	278	459466	39.67	ng #	80
91) Benzo(g,h,i)perylene	19.05	276	465306	39.79	ng #	77

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

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