

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122415\
 Data File : BF084107.D
 Acq On : 25 Dec 2015 4:34
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

Sohil
 12/28/2015 1:10:10 PM

Quant Time: Dec 25 06:19:58 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 24 16:48:39 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	48920	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	183518	20.00	ng	0.00
38) Acenaphthene-d10	9.84	164	87061	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	156198	20.00	ng	0.00
75) Chrysene-d12	13.96	240	117201	20.00	ng	0.00
86) Perylene-d12	15.38	264	93733	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.41	112	225636	78.20	ng	0.00
7) Phenol-d6	6.45	99	262519	73.53	ng	0.00
23) Nitrobenzene-d5	7.37	82	251001	80.06	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	70033	75.43	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	490245	77.05	ng	0.00
78) Terphenyl-d14	12.90	244	469162	72.43	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.34	88	46671	39.46	ng	100
3) Pyridine	3.06	79	123554	42.78	ng	96
4) n-Nitrosodimethylamine	3.00	42	47933	37.67	ng	96
6) Aniline	6.46	93	180490	39.99	ng	# 63
8) 2-Chlorophenol	6.59	128	140709	39.27	ng	90
9) Benzaldehyde	6.35	77	69934	36.62	ng	89
10) Phenol	6.46	94	155247	38.02	ng	# 73
11) bis(2-Chloroethyl)ether	6.53	93	110382	37.48	ng	97
12) 1,3-Dichlorobenzene	6.74	146	148711	38.01	ng	97
13) 1,4-Dichlorobenzene	6.82	146	148633m	37.76	ng	
14) 1,2-Dichlorobenzene	6.97	146	136406	37.60	ng	98
15) Benzyl Alcohol	6.94	79	98875	36.57	ng	# 87
16) 2,2'-oxybis(1-Chloropropan	7.08	45	105839	36.79	ng	66
17) 2-Methylphenol	7.07	107	100267	36.44	ng	# 90
18) Hexachloroethane	7.31	117	55936	38.93	ng	97
19) n-Nitroso-di-n-propylamine	7.22	70	83258	37.86	ng	# 88
20) 3+4-Methylphenols	7.23	107	131833	37.56	ng	# 48
22) Acetophenone	7.22	105	187183	40.82	ng	# 90
24) Nitrobenzene	7.39	77	132730	39.87	ng	# 86
25) Isophorone	7.63	82	231072	39.84	ng	95
26) 2-Nitrophenol	7.71	139	67829	39.81	ng	94
27) 2,4-Dimethylphenol	7.76	122	119967	42.43	ng	93
28) bis(2-Chloroethoxy)methane	7.84	93	127493	37.71	ng	100
29) 2,4-Dichlorophenol	7.95	162	104383	39.75	ng	95
30) 1,2,4-Trichlorobenzene	8.03	180	114071	39.63	ng	97
31) Naphthalene	8.11	128	389759	39.92	ng	98
32) Benzoic acid	7.89	122	53785	39.67	ng	98
33) 4-Chloroaniline	8.16	127	149551	38.49	ng	98
34) Hexachlorobutadiene	8.22	225	63787	38.86	ng	98
35) Caprolactam	8.53	113	30819	42.31	ng	97
36) 4-Chloro-3-methylphenol	8.66	107	121720	41.22	ng	96
37) 2-Methylnaphthalene	8.80	142	235600	38.55	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	109181	39.78	ng	99
40) Hexachlorocyclopentadiene	8.96	237	19450	34.62	ng	96

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42) 2,4,6-Trichlorophenol	9.08	196	70596	39.45	ng	95
43) 2,4,5-Trichlorophenol	9.13	196	75357	42.07	ng #	84
45) 1,1'-Biphenyl	9.26	154	323657	41.15	ng	97
46) 2-Chloronaphthalene	9.29	162	248220	41.51	ng	95
47) 2-Nitroaniline	9.39	65	65813	41.42	ng #	59
48) Acenaphthylene	9.70	152	383794	39.08	ng	100
49) Dimethylphthalate	9.56	163	264244	36.65	ng #	90
50) 2,6-Dinitrotoluene	9.63	165	65633	42.88	ng	95
51) Acenaphthene	9.87	154	224715	39.35	ng	100
52) 3-Nitroaniline	9.80	138	71192	43.34	ng #	73
53) 2,4-Dinitrophenol	9.92	184	19338	42.47	ng	97
54) Dibenzofuran	10.04	168	304727	38.80	ng	93
55) 4-Nitrophenol	9.98	139	32029	38.20	ng	97
56) 2,4-Dinitrotoluene	10.03	165	75385	37.31	ng #	85
57) Fluorene	10.38	166	264172	39.58	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.17	232	52816	41.36	ng	98
59) Diethylphthalate	10.26	149	263546	35.98	ng	95
60) 4-Chlorophenyl-phenylether	10.37	204	108743	35.32	ng #	82
61) 4-Nitroaniline	10.41	138	61540	40.54	ng	95
62) Azobenzene	10.53	77	235185	38.97	ng	91
64) 4,6-Dinitro-2-methylphenol	10.45	198	35012	40.38	ng	80
65) n-Nitrosodiphenylamine	10.50	169	238220	42.78	ng	98
66) 4-Bromophenyl-phenylether	10.87	248	73683	40.72	ng #	82
67) Hexachlorobenzene	10.93	284	75630	38.18	ng #	75
68) Atrazine	11.03	200	67799	40.43	ng	99
69) Pentachlorophenol	11.14	266	24212	41.30	ng	100
70) Phenanthrene	11.35	178	366444	39.57	ng	99
71) Anthracene	11.39	178	369129	38.12	ng	99
72) Carbazole	11.55	167	321606	39.04	ng	99
73) Di-n-butylphthalate	11.88	149	419714	39.55	ng #	99
74) Fluoranthene	12.53	202	389781	39.20	ng	95
76) Benzidine	12.65	184	96598	24.25	ng	100
77) Pyrene	12.76	202	381927	37.77	ng	99
79) Butylbenzylphthalate	13.38	149	164899	39.40	ng #	73
80) Benzo(a)anthracene	13.95	228	291316	40.41	ng	98
81) 3,3'-Dichlorobenzidine	13.91	252	96577	44.24	ng #	96
82) Chrysene	13.99	228	284639	42.82	ng	99
83) Bis(2-ethylhexyl)phthalate	13.94	149	222160	41.08	ng #	97
84) Di-n-octyl phthalate	14.55	149	337209	44.31	ng	99
85) Indeno(1,2,3-cd)pyrene	16.76	276	272167	40.88	ng	100
87) Benzo(b)fluoranthene	14.98	252	279689m	40.18	ng	
88) Benzo(k)fluoranthene	15.00	252	206533m	43.78	ng	
89) Benzo(a)pyrene	15.32	252	227280	39.63	ng	99
90) Dibenzo(a,h)anthracene	16.77	278	223441	39.93	ng #	96
91) Benzo(g,h,i)perylene	17.19	276	228663	40.03	ng	98

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

