

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102615\
 Data File : BF082527.D
 Acq On : 27 Oct 2015 3:11
 Operator : UM/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 27 06:58:29 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 27 02:43:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	101236	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	396894	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	195644	20.00	ng	0.00
63) Phenanthrene-d10	11.26	188	366850	20.00	ng	0.00
75) Chrysene-d12	13.91	240	334579	20.00	ng	0.00
86) Perylene-d12	15.34	264	288524	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.31	112	421375	82.17	ng	0.00
7) Phenol-d6	6.38	99	518644	77.60	ng	0.00
23) Nitrobenzene-d5	7.30	82	480801	78.15	ng	0.00
41) 2,4,6-Tribromophenol	10.57	330	119271	81.41	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	900338	73.18	ng	0.00
78) Terphenyl-d14	12.85	244	838039	72.16	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.23	88	92605	35.98	ng	96
3) Pyridine	2.91	79	251353	41.89	ng	98
4) n-Nitrosodimethylamine	2.89	42	103899	39.38	ng	98
6) Aniline	6.39	93	347584	41.27	ng	96
8) 2-Chlorophenol	6.51	128	249834	41.54	ng	99
9) Benzaldehyde	6.27	77	138022	41.95	ng	98
10) Phenol	6.40	94	278504	36.70	ng	91
11) bis(2-Chloroethyl)ether	6.47	93	230078	38.21	ng	99
12) 1,3-Dichlorobenzene	6.66	146	286952	39.99	ng	98
13) 1,4-Dichlorobenzene	6.74	146	292203	39.58	ng	100
14) 1,2-Dichlorobenzene	6.90	146	256784	37.48	ng	99
15) Benzyl Alcohol	6.89	79	193155	41.52	ng	# 90
16) 2,2'-oxybis(1-Chloropropan	7.02	45	327990	40.23	ng	97
17) 2-Methylphenol	7.01	107	197084	40.86	ng	98
18) Hexachloroethane	7.23	117	100136	39.85	ng	99
19) n-Nitroso-di-n-propylamine	7.15	70	163000	37.43	ng	99
20) 3+4-Methylphenols	7.17	107	244851	39.73	ng	97
22) Acetophenone	7.15	105	339877	39.47	ng	98
24) Nitrobenzene	7.33	77	274074	39.90	ng	97
25) Isophorone	7.57	82	469928	38.45	ng	99
26) 2-Nitrophenol	7.65	139	134507	40.53	ng	98
27) 2,4-Dimethylphenol	7.69	122	223550	41.45	ng	97
28) bis(2-Chloroethoxy)methane	7.78	93	279267	38.96	ng	99
29) 2,4-Dichlorophenol	7.89	162	216795	40.30	ng	98
30) 1,2,4-Trichlorobenzene	7.97	180	237453	39.04	ng	99
31) Naphthalene	8.05	128	689389	38.01	ng	99
32) Benzoic acid	7.85	122	109202m	36.14	ng	
33) 4-Chloroaniline	8.10	127	292364	41.58	ng	96
34) Hexachlorobutadiene	8.15	225	130947	37.22	ng	98
35) Caprolactam	8.49	113	57325	38.34	ng	96
36) 4-Chloro-3-methylphenol	8.59	107	214897	41.08	ng	96
37) 2-Methylnaphthalene	8.73	142	475021	38.77	ng	100
39) 1,2,4,5-Tetrachlorobenzene	8.90	216	237651	40.24	ng	99
40) Hexachlorocyclopentadiene	8.88	237	38419	35.54	ng	100

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42) 2,4,6-Trichlorophenol	9.02	196	144445	39.24	nq	100
43) 2,4,5-Trichlorophenol	9.06	196	157394	40.77	nq	# 92
45) 1,1'-Biphenyl	9.20	154	603069	41.42	nq	99
46) 2-Chloronaphthalene	9.22	162	463791	40.41	nq	99
47) 2-Nitroaniline	9.34	65	138424	41.53	nq	# 55
48) Acenaphthylene	9.63	152	687724	37.42	nq	99
49) Dimethylphthalate	9.51	163	532304	39.10	nq	100
50) 2,6-Dinitrotoluene	9.58	165	125483	42.04	nq	# 83
51) Acenaphthene	9.81	154	403457	38.28	nq	99
52) 3-Nitroaniline	9.75	138	138445	42.74	nq	92
53) 2,4-Dinitrophenol	9.86	184	39776	37.10	nq	# 86
54) Dibenzofuran	9.98	168	572767	37.24	nq	99
55) 4-Nitrophenol	9.93	139	39084m	35.84	nq	
56) 2,4-Dinitrotoluene	9.98	165	146274	40.41	nq	90
57) Fluorene	10.32	166	466685	36.96	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.10	232	120369	40.73	nq	97
59) Diethylphthalate	10.21	149	534915	40.37	nq	100
60) 4-Chlorophenyl-phenylether	10.32	204	226089	38.07	nq	95
61) 4-Nitroaniline	10.37	138	136724	42.15	nq	97
62) Azobenzene	10.48	77	499938	39.07	nq	94
64) 4,6-Dinitro-2-methylphenol	10.40	198	84228	39.14	nq	# 62
65) n-Nitrosodiphenylamine	10.43	169	425957	38.38	nq	99
66) 4-Bromophenyl-phenylether	10.80	248	127903	37.15	nq	94
67) Hexachlorobenzene	10.87	284	149452	40.70	nq	# 91
68) Atrazine	10.97	200	138716	40.37	nq	98
69) Pentachlorophenol	11.07	266	46198	35.34	nq	95
70) Phenanthrene	11.28	178	703220	37.25	nq	100
71) Anthracene	11.34	178	760259	39.99	nq	99
72) Carbazole	11.50	167	648474	38.68	nq	98
73) Di-n-butylphthalate	11.83	149	775923	38.80	nq	# 97
74) Fluoranthene	12.47	202	752260	38.82	nq	99
76) Benzidine	12.61	184	373040	40.63	nq	99
77) Pyrene	12.70	202	766497	37.25	nq	100
79) Butylbenzylphthalate	13.33	149	342364	39.57	nq	# 71
80) Benzo(a)anthracene	13.90	228	687559	38.33	nq	99
81) 3,3'-Dichlorobenzidine	13.86	252	255007	41.00	nq	98
82) Chrysene	13.93	228	619836	36.58	nq	99
83) Bis(2-ethylhexyl)phthalate	13.88	149	435233	37.56	nq	# 94
84) Di-n-octyl phthalate	14.52	149	769645	40.25	nq	99
85) Indeno(1,2,3-cd)pyrene	16.71	276	600443	38.57	nq	99
87) Benzo(b)fluoranthene	14.94	252	583643m	34.06	nq	
88) Benzo(k)fluoranthene	14.97	252	638802m	45.73	nq	
89) Benzo(a)pyrene	15.28	252	573930	38.66	nq	# 96
90) Dibenzo(a,h)anthracene	16.71	278	500933	40.33	nq	# 96
91) Benzo(g,h,i)perylene	17.12	276	485277	39.49	nq	98

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

