

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
 Data File : BF081286.D
 Acq On : 29 Aug 2015 22:59
 Operator : UM/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

MMDadoda
 8/31/2015 1:59:13 PM

Quant Time: Aug 31 02:19:43 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 31 00:56:51 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.48	152	65525	20.00	ng	0.00
21) Naphthalene-d8	7.77	136	266482	20.00	ng	0.00
38) Acenaphthene-d10	9.52	164	128106	20.00	ng	0.00
63) Phenanthrene-d10	10.98	188	241564	20.00	ng	0.00
75) Chrysene-d12	13.60	240	192560	20.00	ng	0.00
86) Perylene-d12	14.92	264	135183	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.04	112	367697	84.40	ng	0.01
7) Phenol-d6	6.15	99	456154	80.25	ng	0.01
23) Nitrobenzene-d5	7.06	82	431513	73.97	ng	0.00
41) 2,4,6-Tribromophenol	10.31	330	84191	75.82	ng	0.00
44) 2-Fluorobiphenyl	8.86	172	687611	70.33	ng	0.00
78) Terphenyl-d14	12.57	244	727266	80.97	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.85	88	83352	40.60	ng	# 89
3) Pyridine	2.43	79	255779	44.14	ng	# 82
4) n-Nitrosodimethylamine	2.41	42	127613	39.56	ng	# 77
6) Aniline	6.15	93	318240	38.51	ng	# 71
8) 2-Chlorophenol	6.26	128	188857	39.61	ng	94
9) Benzaldehyde	6.02	77	148118	40.42	ng	97
10) Phenol	6.16	94	297364	44.30	ng	# 76
11) bis(2-Chloroethyl)ether	6.24	93	217478	42.22	ng	97
12) 1,3-Dichlorobenzene	6.42	146	217516	42.45	ng	99
13) 1,4-Dichlorobenzene	6.50	146	221529	43.66	ng	98
14) 1,2-Dichlorobenzene	6.65	146	164616	34.67	ng	97
15) Benzyl Alcohol	6.65	79	195497	42.51	ng	91
16) 2,2'-oxybis(1-Chloropropan	6.79	45	393971	38.92	ng	96
17) 2-Methylphenol	6.77	107	171508	41.92	ng	96
18) Hexachloroethane	6.99	117	84607	41.27	ng	99
19) n-Nitroso-di-n-propylamine	6.93	70	156563	39.76	ng	90
20) 3+4-Methylphenols	6.94	107	218934	42.16	ng	# 40
22) Acetophenone	6.91	105	286385	40.94	ng	# 85
24) Nitrobenzene	7.09	77	268507	44.02	ng	94
25) Isophorone	7.33	82	441848	40.62	ng	99
26) 2-Nitrophenol	7.39	139	94325	37.19	ng	89
27) 2,4-Dimethylphenol	7.45	122	169584	37.79	ng	95
28) bis(2-Chloroethoxy)methane	7.55	93	261199	40.64	ng	# 98
29) 2,4-Dichlorophenol	7.65	162	158660	42.00	ng	92
30) 1,2,4-Trichlorobenzene	7.71	180	162528	38.71	ng	97
31) Naphthalene	7.79	128	544118	36.63	ng	99
32) Benzoic acid	7.62	122	119223	47.23	ng	90
33) 4-Chloroaniline	7.86	127	246913	39.40	ng	# 82
34) Hexachlorobutadiene	7.92	225	99603	41.95	ng	96
35) Caprolactam	8.26	113	54782	41.49	ng	# 44
36) 4-Chloro-3-methylphenol	8.35	107	167071	37.11	ng	99
37) 2-Methylnaphthalene	8.49	142	360787m	38.45	ng	
39) 1,2,4,5-Tetrachlorobenzene	8.65	216	142517	34.48	ng	# 97
40) Hexachlorocyclopentadiene	8.64	237	81071	37.90	ng	97

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42) 2,4,6-Trichlorophenol	8.78	196	109816	37.61	ng	97
43) 2,4,5-Trichlorophenol	8.81	196	111379	38.17	ng #	90
45) 1,1'-Biphenyl	8.95	154	369922	32.78	ng	95
46) 2-Chloronaphthalene	8.97	162	330222	36.43	ng	94
47) 2-Nitroaniline	9.07	65	124974	33.69	ng	96
48) Acenaphthylene	9.38	152	595125	36.70	ng	99
49) Dimethylphthalate	9.27	163	363129	34.42	ng	99
50) 2,6-Dinitrotoluene	9.33	165	87235	38.51	ng #	84
51) Acenaphthene	9.55	154	341596	35.19	ng	100
52) 3-Nitroaniline	9.49	138	96556	34.06	ng	99
53) 2,4-Dinitrophenol	9.60	184	34109	32.29	ng #	64
54) Dibenzofuran	9.73	168	469792	36.11	ng	99
55) 4-Nitrophenol	9.67	139	73290	36.81	ng #	71
56) 2,4-Dinitrotoluene	9.73	165	115947	38.62	ng #	88
57) Fluorene	10.06	166	319421	31.92	ng	99
58) 2,3,4,6-Tetrachlorophenol	9.85	232	80393	35.59	ng	96
59) Diethylphthalate	9.97	149	437421	39.17	ng #	94
60) 4-Chlorophenyl-phenylether	10.06	204	151361	35.74	ng	96
61) 4-Nitroaniline	10.10	138	100956	34.85	ng	93
62) Azobenzene	10.22	77	439012	34.12	ng	96
64) 4,6-Dinitro-2-methylphenol	10.14	198	56498	39.16	ng #	53
65) n-Nitrosodiphenylamine	10.18	169	300550	34.85	ng	99
66) 4-Bromophenyl-phenylether	10.55	248	97773	41.20	ng #	94
67) Hexachlorobenzene	10.61	284	103728	43.16	ng	93
68) Atrazine	10.72	200	89000	37.04	ng	99
69) Pentachlorophenol	10.80	266	20770	14.40	ng	98
70) Phenanthrene	11.01	178	474016	33.11	ng	99
71) Anthracene	11.06	178	569304	40.87	ng	100
72) Carbazole	11.22	167	523479	39.03	ng	99
73) Di-n-butylphthalate	11.57	149	639558	39.41	ng	98
74) Fluoranthene	12.18	202	540450	34.98	ng	94
76) Benzidine	12.32	184	309453	42.54	ng	99
77) Pyrene	12.41	202	614294	39.24	ng	99
79) Butylbenzylphthalate	13.05	149	298359	44.34	ng #	80
80) Benzo(a)anthracene	13.59	228	495614	38.38	ng	100
81) 3,3'-Dichlorobenzidine	13.57	252	158025	37.78	ng	99
82) Chrysene	13.62	228	373257	32.07	ng	100
83) Bis(2-ethylhexyl)phthalate	13.61	149	370838	43.03	ng #	97
84) Di-n-octyl phthalate	14.22	149	651046	49.70	ng	97
85) Indeno(1,2,3-cd)pyrene	16.07	276	333645	40.83	ng #	94
87) Benzo(b)fluoranthene	14.58	252	442213m	46.25	ng	
88) Benzo(k)fluoranthene	14.61	252	300884m	33.77	ng	
89) Benzo(a)pyrene	14.87	252	321792	38.62	ng #	95
90) Dibenzo(a,h)anthracene	16.08	278	276147	42.54	ng #	95
91) Benzo(g,h,i)perylene	16.41	276	271376	41.20	ng #	94

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

