

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061116\
 Data File : BF087921.D
 Acq On : 12 Jun 2016 00:28
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

umangi
 6/13/2016 7:44:35 PM

Quant Time: Jun 12 05:56:04 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 07 18:51:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	105590	20.00	ng	-0.05
21) Naphthalene-d8	8.09	136	430089	20.00	ng	-0.03
38) Acenaphthene-d10	9.85	164	252866	20.00	ng	-0.03
63) Phenanthrene-d10	11.33	188	444940	20.00	ng	-0.03
75) Chrysene-d12	13.97	240	348338	20.00	ng	-0.02
86) Perylene-d12	15.37	264	278234	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	507971	82.24	ng	-0.05
7) Phenol-d6	6.43	99	714880	95.50	ng	-0.05
23) Nitrobenzene-d5	7.37	82	563093	76.45	ng	-0.05
41) 2,4,6-Tribromophenol	10.64	330	184805	69.22	ng	-0.03
44) 2-Fluorobiphenyl	9.18	172	1192402	73.32	ng	-0.03
78) Terphenyl-d14	12.91	244	966284	63.12	ng	-0.03

Target Compounds						Qvalue
2) 1,4-Dioxane	2.34	88	126748	42.23	ng	# 40
3) Pyridine	3.04	79	384372	54.24	ng	97
4) n-Nitrosodimethylamine	2.99	42	139256	46.40	ng	89
6) Aniline	6.47	93	494390	46.40	ng	97
8) 2-Chlorophenol	6.58	128	281107	38.45	ng	98
9) Benzaldehyde	6.34	77	194786	45.98	ng	91
10) Phenol	6.46	94	418042	54.62	ng	# 69
11) bis(2-Chloroethyl)ether	6.55	93	271500	41.36	ng	# 81
12) 1,3-Dichlorobenzene	6.74	146	302448	38.56	ng	98
13) 1,4-Dichlorobenzene	6.82	146	315232	39.89	ng	98
14) 1,2-Dichlorobenzene	6.97	146	263144m	36.62	ng	
15) Benzyl Alcohol	6.95	79	209513	35.78	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.10	45	352493	39.75	ng	89
17) 2-Methylphenol	7.06	107	240500	40.57	ng	98
18) Hexachloroethane	7.31	117	108443	38.68	ng	# 83
19) n-Nitroso-di-n-propylamine	7.23	70	206386	43.10	ng	# 84
20) 3+4-Methylphenols	7.22	107	256762	37.12	ng	# 80
22) Acetophenone	7.22	105	365533	38.03	ng	# 90
24) Nitrobenzene	7.39	77	322449	45.20	ng	95
25) Isophorone	7.63	82	543228	39.82	ng	100
26) 2-Nitrophenol	7.71	139	163851	42.87	ng	# 70
27) 2,4-Dimethylphenol	7.75	122	269583	38.31	ng	94
28) bis(2-Chloroethoxy)methane	7.85	93	361081	42.67	ng	98
29) 2,4-Dichlorophenol	7.95	162	248175	42.24	ng	96
30) 1,2,4-Trichlorobenzene	8.03	180	242097	37.84	ng	99
31) Naphthalene	8.11	128	797908	37.49	ng	100
32) Benzoic acid	7.88	122	169150	43.65	ng	91
33) 4-Chloroaniline	8.17	127	399361	42.02	ng	# 91
34) Hexachlorobutadiene	8.24	225	134953	40.00	ng	99
35) Caprolactam	8.55	113	75844	38.97	ng	91
36) 4-Chloro-3-methylphenol	8.65	107	249967	39.78	ng	100
37) 2-Methylnaphthalene	8.81	142	608547	43.38	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	228722	31.46	ng	# 1
40) Hexachlorocyclopentadiene	8.96	237	124365	30.49	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.08	196	192443	38.06	nq	98
43) 2,4,5-Trichlorophenol	9.13	196	192643	36.62	nq	# 94
45) 1,1'-Biphenyl	9.27	154	721703	36.91	nq	99
46) 2-Chloronaphthalene	9.29	162	590927	36.11	nq	100
47) 2-Nitroaniline	9.39	65	201747	41.80	nq	# 76
48) Acenaphthylene	9.71	152	992374	38.13	nq	100
49) Dimethylphthalate	9.58	163	635040	34.67	nq	99
50) 2,6-Dinitrotoluene	9.63	165	146645	34.96	nq	# 60
51) Acenaphthene	9.88	154	625507	38.52	nq	99
52) 3-Nitroaniline	9.80	138	190491	39.35	nq	91
53) 2,4-Dinitrophenol	9.91	184	87078	43.57	nq	# 85
54) Dibenzofuran	10.06	168	899386	40.22	nq	97
55) 4-Nitrophenol	9.96	139	153325	40.81	nq	# 84
56) 2,4-Dinitrotoluene	10.03	165	197759	36.68	nq	# 63
57) Fluorene	10.40	166	617494	40.37	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.17	232	160723	38.87	nq	# 52
59) Diethylphthalate	10.27	149	652443	35.19	nq	99
60) 4-Chlorophenyl-phenylether	10.39	204	250159	33.64	nq	99
61) 4-Nitroaniline	10.41	138	191823	41.78	nq	97
62) Azobenzene	10.55	77	692454	37.77	nq	98
64) 4,6-Dinitro-2-methylphenol	10.44	198	114489	41.52	nq	80
65) n-Nitrosodiphenylamine	10.51	169	603152	40.85	nq	100
66) 4-Bromophenyl-phenylether	10.88	248	189748	39.75	nq	98
67) Hexachlorobenzene	10.94	284	182244	36.74	nq	95
68) Atrazine	11.04	200	183407	41.02	nq	94
69) Pentachlorophenol	11.13	266	111107	42.50	nq	97
70) Phenanthrene	11.35	178	884253	37.87	nq	100
71) Anthracene	11.41	178	913279	38.78	nq	100
72) Carbazole	11.55	167	891828	40.47	nq	100
73) Di-n-butylphthalate	11.90	149	986488	35.36	nq	100
74) Fluoranthene	12.54	202	900395	36.54	nq	100
76) Benzidine	12.66	184	461283	47.83	nq	99
77) Pyrene	12.77	202	926883	35.86	nq	100
79) Butylbenzylphthalate	13.39	149	470938	41.21	nq	97
80) Benzo(a)anthracene	13.95	228	697394	35.13	nq	99
81) 3,3'-Dichlorobenzidine	13.92	252	235777	44.17	nq	# 97
82) Chrysene	13.99	228	673864	36.08	nq	100
83) Bis(2-ethylhexyl)phthalate	13.95	149	502015	36.09	nq	# 98
84) Di-n-octyl phthalate	14.56	149	917729	40.60	nq	# 100
85) Indeno(1,2,3-cd)pyrene	16.74	276	638457	36.80	nq	# 100
87) Benzo(b)fluoranthene	14.97	252	634027m	32.69	nq	
88) Benzo(k)fluoranthene	15.01	252	671148m	45.88	nq	
89) Benzo(a)pyrene	15.31	252	619233	39.47	nq	99
90) Dibenzo(a,h)anthracene	16.77	278	524676	36.00	nq	98
91) Benzo(g,h,i)perylene	17.15	276	543656	36.27	nq	98

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

