

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062016\
 Data File : BF088189.D
 Acq On : 20 Jun 2016 23:48
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

sohil
 6/21/2016 4:22:52 PM

Quant Time: Jun 21 01:47:46 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 17:39:32 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	81430	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	336136	20.00	ng	0.00
38) Acenaphthene-d10	9.70	164	156513	20.00	ng	0.00
63) Phenanthrene-d10	11.18	188	288484	20.00	ng	0.00
75) Chrysene-d12	13.81	240	215258	20.00	ng	0.00
86) Perylene-d12	15.18	264	186595	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.26	112	385202	80.87	ng	0.00
7) Phenol-d6	6.32	99	431767	74.79	ng	-0.01
23) Nitrobenzene-d5	7.24	82	443377	77.02	ng	0.00
41) 2,4,6-Tribromophenol	10.49	330	106658	64.54	ng	0.00
44) 2-Fluorobiphenyl	9.03	172	806975	80.72	ng	-0.01
78) Terphenyl-d14	12.77	244	765021	80.87	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.12	88	101607	43.90	ng	# 43
3) Pyridine	2.79	79	207021	37.88	ng	91
4) n-Nitrosodimethylamine	2.74	42	86798	37.50	ng	79
6) Aniline	6.33	93	282470	34.38	ng	# 67
8) 2-Chlorophenol	6.46	128	232802	41.29	ng	88
9) Benzaldehyde	6.22	77	132064	34.94	ng	99
10) Phenol	6.34	94	246104	41.07	ng	91
11) bis(2-Chloroethyl)ether	6.41	93	201300	39.77	ng	89
12) 1,3-Dichlorobenzene	6.60	146	237181	39.21	ng	98
13) 1,4-Dichlorobenzene	6.68	146	248726	40.82	ng	99
14) 1,2-Dichlorobenzene	6.84	146	210414	37.97	ng	99
15) Benzyl Alcohol	6.82	79	148879	32.97	ng	95
16) 2,2'-oxybis(1-Chloropropan	6.96	45	235135	34.38	ng	78
17) 2-Methylphenol	6.95	107	179766	39.32	ng	99
18) Hexachloroethane	7.18	117	89060	41.19	ng	90
19) n-Nitroso-di-n-propylamine	7.10	70	147585	39.96	ng	# 86
20) 3+4-Methylphenols	7.11	107	228008	42.74	ng	# 74
22) Acetophenone	7.08	105	292553	38.95	ng	# 96
24) Nitrobenzene	7.26	77	201731	36.18	ng	# 87
25) Isophorone	7.50	82	394818	37.03	ng	98
26) 2-Nitrophenol	7.58	139	122133	40.89	ng	# 88
27) 2,4-Dimethylphenol	7.63	122	193809	35.24	ng	99
28) bis(2-Chloroethoxy)methane	7.71	93	247763	37.47	ng	# 96
29) 2,4-Dichlorophenol	7.83	162	183024	39.86	ng	93
30) 1,2,4-Trichlorobenzene	7.90	180	187768	37.55	ng	98
31) Naphthalene	7.98	128	641717	38.58	ng	99
32) Benzoic acid	7.76	122	66457	21.95	ng	96
33) 4-Chloroaniline	8.03	127	271973	36.62	ng	97
34) Hexachlorobutadiene	8.09	225	98128	37.21	ng	97
35) Caprolactam	8.41	113	58995	38.79	ng	# 66
36) 4-Chloro-3-methylphenol	8.54	107	187391	38.16	ng	# 82
37) 2-Methylnaphthalene	8.66	142	406471	37.07	ng	# 96
39) 1,2,4,5-Tetrachlorobenzene	8.83	216	178370	43.20	ng	# 1
40) Hexachlorocyclopentadiene	8.82	237	85915	34.03	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.96	196	109045	34.84	nq	99
43) 2,4,5-Trichlorophenol	8.99	196	117367	36.04	nq	# 78
45) 1,1'-Biphenyl	9.13	154	489183	41.10	nq	98
46) 2-Chloronaphthalene	9.15	162	393699	38.87	nq	98
47) 2-Nitroaniline	9.26	65	107638	36.03	nq	92
48) Acenaphthylene	9.56	152	582541	36.16	nq	100
49) Dimethylphthalate	9.44	163	448850	39.59	nq	100
50) 2,6-Dinitrotoluene	9.51	165	103199	39.75	nq	# 81
51) Acenaphthene	9.74	154	362812	36.10	nq	98
52) 3-Nitroaniline	9.67	138	106010	35.38	nq	95
53) 2,4-Dinitrophenol	9.78	184	45297	37.73	nq	98
54) Dibenzofuran	9.91	168	494727	35.74	nq	# 91
55) 4-Nitrophenol	9.85	139	90709	39.01	nq	93
56) 2,4-Dinitrotoluene	9.91	165	123596	37.04	nq	# 71
57) Fluorene	10.25	166	421331	45.05	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.03	232	95429	37.29	nq	# 59
59) Diethylphthalate	10.14	149	482089	42.01	nq	99
60) 4-Chlorophenyl-phenylether	10.24	204	169141	36.75	nq	89
61) 4-Nitroaniline	10.28	138	125270	44.08	nq	99
62) Azobenzene	10.40	77	429779	37.87	nq	94
64) 4,6-Dinitro-2-methylphenol	10.32	198	51262	29.34	nq	# 58
65) n-Nitrosodiphenylamine	10.36	169	374428	39.11	nq	99
66) 4-Bromophenyl-phenylether	10.73	248	120287	38.87	nq	# 92
67) Hexachlorobenzene	10.80	284	122844	38.20	nq	# 71
68) Atrazine	10.89	200	106784	36.84	nq	91
69) Pentachlorophenol	10.99	266	59862	35.32	nq	98
70) Phenanthrene	11.20	178	599492	39.60	nq	99
71) Anthracene	11.26	178	616535	40.38	nq	99
72) Carbazole	11.42	167	575546	40.28	nq	100
73) Di-n-butylphthalate	11.75	149	691481	38.23	nq	99
74) Fluoranthene	12.39	202	638767	39.98	nq	97
76) Benzidine	12.51	184	252181	42.31	nq	99
77) Pyrene	12.62	202	635904	39.81	nq	98
79) Butylbenzylphthalate	13.23	149	292731	41.45	nq	88
80) Benzo(a)anthracene	13.79	228	451562	36.81	nq	99
81) 3,3'-Dichlorobenzidine	13.76	252	140137	42.48	nq	# 95
82) Chrysene	13.84	228	496662	43.04	nq	98
83) Bis(2-ethylhexyl)phthalate	13.79	149	360737	41.96	nq	99
84) Di-n-octyl phthalate	14.40	149	681888	48.81	nq	# 100
85) Indeno(1,2,3-cd)pyrene	16.47	276	419484	39.12	nq	# 100
87) Benzo(b)fluoranthene	14.80	252	417581m	32.11	nq	
88) Benzo(k)fluoranthene	14.83	252	445369m	45.40	nq	
89) Benzo(a)pyrene	15.13	252	415106	39.45	nq	# 93
90) Dibenzo(a,h)anthracene	16.47	278	348135	35.62	nq	96
91) Benzo(g,h,i)perylene	16.85	276	367588	36.57	nq	97

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

