

Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\
 Data File : BF089465.D
 Acq On : 31 Jul 2016 10:53
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

sohil
 8/1/2016 7:18:06 PM

Quant Time: Aug 01 12:37:30 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 04:11:30 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.49	152	49134	20.00	ng	-0.01
21) Naphthalene-d8	7.78	136	210379	20.00	ng	-0.01
38) Acenaphthene-d10	9.53	164	100190	20.00	ng	0.00
63) Phenanthrene-d10	10.99	188	195292	20.00	ng	-0.01
75) Chrysene-d12	13.61	240	134961	20.00	ng	-0.01
86) Perylene-d12	14.95	264	110458	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.05	112	248285	80.69	ng	-0.01
7) Phenol-d6	6.16	99	323573	79.57	ng	-0.01
23) Nitrobenzene-d5	7.07	82	313228	87.88	ng	-0.01
41) 2,4,6-Tribromophenol	10.32	330	67609	81.48	ng	0.00
44) 2-Fluorobiphenyl	8.87	172	575073	84.24	ng	0.00
78) Terphenyl-d14	12.58	244	523902	90.85	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.87	88	57831	41.62	ng	97
3) Pyridine	2.46	79	141753	47.08	ng	99
4) n-Nitrosodimethylamine	2.43	42	49657	41.34	ng	97
6) Aniline	6.16	93	207439	47.24	ng	96
8) 2-Chlorophenol	6.27	128	139102	39.96	ng	86
9) Benzaldehyde	6.03	77	70317m	35.41	ng	
10) Phenol	6.17	94	196862	43.87	ng	90
11) bis(2-Chloroethyl)ether	6.24	93	147530	38.69	ng	90
12) 1,3-Dichlorobenzene	6.43	146	142069m	39.02	ng	
13) 1,4-Dichlorobenzene	6.51	146	158729	39.50	ng	96
14) 1,2-Dichlorobenzene	6.66	146	154375	41.01	ng	95
15) Benzyl Alcohol	6.66	79	108486	42.99	ng	97
16) 2,2'-oxybis(1-Chloropropan	6.79	45	165538	39.80	ng	99
17) 2-Methylphenol	6.79	107	100887	37.50	ng	# 90
18) Hexachloroethane	7.00	117	61886	40.17	ng	# 82
19) n-Nitroso-di-n-propylamine	6.94	70	106572	43.46	ng	99
20) 3+4-Methylphenols	6.95	107	141792	41.04	ng	# 85
22) Acetophenone	6.92	105	216866	43.30	ng	# 98
24) Nitrobenzene	7.10	77	148185	36.53	ng	93
25) Isophorone	7.34	82	285084	39.71	ng	99
26) 2-Nitrophenol	7.40	139	73930	40.76	ng	# 79
27) 2,4-Dimethylphenol	7.46	122	118426	40.99	ng	99
28) bis(2-Chloroethoxy)methane	7.55	93	180123	45.32	ng	99
29) 2,4-Dichlorophenol	7.66	162	118556	45.03	ng	98
30) 1,2,4-Trichlorobenzene	7.72	180	122155	40.67	ng	# 95
31) Naphthalene	7.80	128	432881	41.55	ng	100
32) Benzoic acid	7.63	122	86243	42.84	ng	91
33) 4-Chloroaniline	7.87	127	170409	42.97	ng	98
34) Hexachlorobutadiene	7.93	225	67941	39.49	ng	97
35) Caprolactam	8.25	113	34566	43.50	ng	93
36) 4-Chloro-3-methylphenol	8.36	107	137307	43.88	ng	98
37) 2-Methylnaphthalene	8.49	142	256212	40.81	ng	96
39) 1,2,4,5-Tetrachlorobenzene	8.66	216	143702	40.99	ng	99
40) Hexachlorocyclopentadiene	8.65	237	32680	37.97	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.79	196	71952	43.16	nq	99
43) 2,4,5-Trichlorophenol	8.83	196	86205	40.83	nq	100
45) 1,1'-Biphenyl	8.96	154	353314	41.42	nq	98
46) 2-Chloronaphthalene	8.98	162	275689	43.03	nq	96
47) 2-Nitroaniline	9.08	65	88141	46.46	nq	# 68
48) Acenaphthylene	9.38	152	397074	39.35	nq	99
49) Dimethylphthalate	9.28	163	305818	40.11	nq	99
50) 2,6-Dinitrotoluene	9.34	165	68004	43.41	nq	97
51) Acenaphthene	9.55	154	232509	39.46	nq	99
52) 3-Nitroaniline	9.51	138	77559	46.91	nq	97
53) 2,4-Dinitrophenol	9.62	184	20710	40.10	nq	91
54) Dibenzofuran	9.74	168	321196	40.00	nq	95
55) 4-Nitrophenol	9.69	139	31669m	42.47	nq	
56) 2,4-Dinitrotoluene	9.74	165	87244	42.30	nq	92
57) Fluorene	10.07	166	264165	37.69	nq	99
58) 2,3,4,6-Tetrachlorophenol	9.86	232	64819	45.98	nq	97
59) Diethylphthalate	9.96	149	291288	37.68	nq	99
60) 4-Chlorophenyl-phenylether	10.07	204	126059	39.21	nq	# 81
61) 4-Nitroaniline	10.11	138	75635	44.38	nq	98
62) Azobenzene	10.23	77	329253	41.36	nq	94
64) 4,6-Dinitro-2-methylphenol	10.15	198	48095	41.45	nq	93
65) n-Nitrosodiphenylamine	10.19	169	255591	41.94	nq	99
66) 4-Bromophenyl-phenylether	10.56	248	73298	39.31	nq	# 90
67) Hexachlorobenzene	10.62	284	80407	42.08	nq	# 80
68) Atrazine	10.73	200	80902	43.22	nq	98
69) Pentachlorophenol	10.82	266	34793	40.80	nq	99
70) Phenanthrene	11.02	178	385748	38.30	nq	99
71) Anthracene	11.07	178	456959	40.81	nq	99
72) Carbazole	11.23	167	386954	41.05	nq	98
73) Di-n-butylphthalate	11.58	149	484155	38.28	nq	99
74) Fluoranthene	12.19	202	412538	38.94	nq	95
76) Benzidine	12.33	184	180135	36.12	nq	98
77) Pyrene	12.42	202	445822	43.59	nq	100
79) Butylbenzylphthalate	13.06	149	203964	43.86	nq	# 82
80) Benzo(a)anthracene	13.60	228	323920	42.53	nq	100
81) 3,3'-Dichlorobenzidine	13.58	252	118610	43.28	nq	# 97
82) Chrysene	13.65	228	347107	42.81	nq	99
83) Bis(2-ethylhexyl)phthalate	13.62	149	286848	42.67	nq	# 97
84) Di-n-octyl phthalate	14.23	149	465556	44.38	nq	98
85) Indeno(1,2,3-cd)pyrene	16.11	276	216885	37.63	nq	99
87) Benzo(b)fluoranthene	14.61	252	325230m	47.41	nq	
88) Benzo(k)fluoranthene	14.63	252	227890m	36.13	nq	
89) Benzo(a)pyrene	14.89	252	242894	39.47	nq	99
90) Dibenzo(a,h)anthracene	16.13	278	191093	39.42	nq	96
91) Benzo(g,h,i)perylene	16.46	276	187382	39.32	nq	99

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

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