

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072716\
 Data File : BF089327.D
 Acq On : 27 Jul 2016 13:32
 Operator : UM/SJ
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations

sohil
 7/28/2016 4:55:49 PM

Quant Time: Jul 27 15:29:12 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 27 15:17:12 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.55	152	45059	20.00	ng	0.00
21) Naphthalene-d8	7.84	136	188515	20.00	ng	0.01
38) Acenaphthene-d10	9.58	164	89549	20.00	ng	0.00
63) Phenanthrene-d10	11.05	188	171797	20.00	ng	0.00
75) Chrysene-d12	13.67	240	113970	20.00	ng	0.00
86) Perylene-d12	15.01	264	93602	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.12	112	345485	122.43	ng	0.01
7) Phenol-d6	6.20	99	437796	117.39	ng	0.00
23) Nitrobenzene-d5	7.13	82	400291	125.33	ng	0.01
41) 2,4,6-Tribromophenol	10.36	330	89788	121.06	ng	0.00
44) 2-Fluorobiphenyl	8.91	172	677238	110.99	ng	0.00
78) Terphenyl-d14	12.63	244	621329	127.59	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.95	88	74783	51.34	ng	99
3) Pyridine	2.56	79	178892	64.79	ng	98
4) n-Nitrosodimethylamine	2.54	42	72079	72.40	ng	96
6) Aniline	6.22	93	250870	62.30	ng	# 86
8) 2-Chlorophenol	6.33	128	192202	60.21	ng	93
9) Benzaldehyde	6.09	77	108225	59.43	ng	97
10) Phenol	6.23	94	244765	59.48	ng	92
11) bis(2-Chloroethyl)ether	6.30	93	196531	56.21	ng	95
12) 1,3-Dichlorobenzene	6.48	146	192251	57.58	ng	97
13) 1,4-Dichlorobenzene	6.56	146	209130	56.76	ng	98
14) 1,2-Dichlorobenzene	6.72	146	202046	58.53	ng	96
15) Benzyl Alcohol	6.71	79	141618	61.19	ng	95
16) 2,2'-oxybis(1-Chloropropan	6.84	45	225287	59.07	ng	81
17) 2-Methylphenol	6.83	107	146709	59.46	ng	96
18) Hexachloroethane	7.05	117	80305	56.84	ng	98
19) n-Nitroso-di-n-propylamine	6.98	70	135112	60.08	ng	# 89
20) 3+4-Methylphenols	6.99	107	188908	59.62	ng	# 85
22) Acetophenone	6.97	105	270363	60.24	ng	# 94
24) Nitrobenzene	7.14	77	213558	58.75	ng	91
25) Isophorone	7.39	82	388743	60.42	ng	97
26) 2-Nitrophenol	7.46	139	101860	62.68	ng	99
27) 2,4-Dimethylphenol	7.52	122	157738	60.93	ng	94
28) bis(2-Chloroethoxy)methane	7.61	93	212548	59.68	ng	98
29) 2,4-Dichlorophenol	7.70	162	147274	62.43	ng	92
30) 1,2,4-Trichlorobenzene	7.78	180	171170	63.59	ng	96
31) Naphthalene	7.85	128	525922	56.34	ng	99
32) Benzoic acid	7.70	122	116653	64.66	ng	97
33) 4-Chloroaniline	7.92	127	222296	62.55	ng	97
34) Hexachlorobutadiene	7.98	225	90190	58.50	ng	99
35) Caprolactam	8.32	113	45744	64.24	ng	97
36) 4-Chloro-3-methylphenol	8.42	107	177083	63.16	ng	91
37) 2-Methylnaphthalene	8.55	142	356857	63.44	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.72	216	187959	59.98	ng	99
40) Hexachlorocyclopentadiene	8.70	237	51544	72.73	ng	99

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42) 2,4,6-Trichlorophenol	8.83	196	96766	64.95	nq	99
43) 2,4,5-Trichlorophenol	8.88	196	116859	61.92	nq	97
45) 1,1'-Biphenyl	9.02	154	459384	60.25	nq	95
46) 2-Chloronaphthalene	9.03	162	333158	58.18	nq	98
47) 2-Nitroaniline	9.14	65	109151	64.37	nq	91
48) Acenaphthylene	9.44	152	506352	56.14	nq	99
49) Dimethylphthalate	9.32	163	393619	57.75	nq	98
50) 2,6-Dinitrotoluene	9.38	165	82624	59.00	nq	# 59
51) Acenaphthene	9.61	154	293560	55.74	nq	99
52) 3-Nitroaniline	9.55	138	88677	60.01	nq	# 80
53) 2,4-Dinitrophenol	9.67	184	36384	78.11	nq	98
54) Dibenzofuran	9.78	168	406267	56.61	nq	97
55) 4-Nitrophenol	9.74	139	44846m	46.16	nq	
56) 2,4-Dinitrotoluene	9.79	165	117851	63.93	nq	# 69
57) Fluorene	10.12	166	357884	57.12	nq	99
58) 2,3,4,6-Tetrachlorophenol	9.91	232	76608	60.80	nq	# 96
59) Diethylphthalate	10.02	149	435969	63.10	nq	96
60) 4-Chlorophenyl-phenylether	10.12	204	174840	60.84	nq	93
61) 4-Nitroaniline	10.17	138	90621	59.49	nq	97
62) Azobenzene	10.28	77	426788	59.98	nq	88
64) 4,6-Dinitro-2-methylphenol	10.20	198	62007	64.15	nq	85
65) n-Nitrosodiphenylamine	10.25	169	327481	61.09	nq	99
66) 4-Bromophenyl-phenylether	10.60	248	97852	59.66	nq	# 81
67) Hexachlorobenzene	10.67	284	102376	60.91	nq	# 72
68) Atrazine	10.78	200	87666	53.23	nq	96
69) Pentachlorophenol	10.87	266	48626	69.44	nq	100
70) Phenanthrene	11.07	178	508585	57.40	nq	100
71) Anthracene	11.13	178	587253	59.61	nq	99
72) Carbazole	11.29	167	501168	60.44	nq	99
73) Di-n-butylphthalate	11.63	149	677798	60.92	nq	# 98
74) Fluoranthene	12.25	202	524174	56.24	nq	96
76) Benzidine	12.39	184	267562	63.54	nq	98
77) Pyrene	12.48	202	565680	65.49	nq	99
79) Butylbenzylphthalate	13.11	149	246358	62.73	nq	87
80) Benzo(a)anthracene	13.66	228	404851	62.95	nq	99
81) 3,3'-Dichlorobenzidine	13.63	252	149538	64.61	nq	# 96
82) Chrysene	13.70	228	444814	64.97	nq	99
83) Bis(2-ethylhexyl)phthalate	13.68	149	368706	64.95	nq	# 94
84) Di-n-octyl phthalate	14.29	149	599413	67.67	nq	99
85) Indeno(1,2,3-cd)pyrene	16.21	276	303267	62.30	nq	99
87) Benzo(b)fluoranthene	14.66	252	423249m	72.81	nq	
88) Benzo(k)fluoranthene	14.69	252	262316m	49.24	nq	
89) Benzo(a)pyrene	14.96	252	320591	61.48	nq	# 95
90) Dibenzo(a,h)anthracene	16.22	278	252618	61.50	nq	98
91) Benzo(g,h,i)perylene	16.56	276	243338	60.26	nq	98

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

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