

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061115\
 Data File : BF079891.D
 Acq On : 11 Jun 2015 13:09
 Operator : TP/IZ
 Sample : SSTDICC025
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC025

Manual Integrations
 APPROVED

apatel
 6/13/2015 8:43:09 AM

Quant Time: Jun 12 08:49:40 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF061115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 12 00:59:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	48484	20.00	ng	0.00
21) Naphthalene-d8	8.68	136	191335	20.00	ng	0.00
38) Acenaphthene-d10	10.46	164	93413	20.00	ng	0.00
63) Phenanthrene-d10	11.99	188	196130	20.00	ng	0.01
75) Chrysene-d12	15.17	240	207804m	20.00	ng	0.17
86) Perylene-d12	17.19	264	194848m	20.00	ng	0.26

System Monitoring Compounds

5) 2-Fluorophenol	6.00	112	131910	45.59	ng	0.00
7) Phenol-d6	6.99	99	178496	47.15	ng	0.00
23) Nitrobenzene-d5	7.95	82	142497	44.30	ng	0.00
41) 2,4,6-Tribromophenol	11.26	330	42302	50.74	ng	0.00
44) 2-Fluorobiphenyl	9.76	172	354632	53.50	ng	0.00
78) Terphenyl-d14	13.78	244	471262	51.53	ng	0.09

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.34	88	30013	22.68	ng	94
3) Pyridine	4.14	79	68967	19.93	ng	99
4) n-Nitrosodimethylamine	4.03	42	39119	22.65	ng	94
6) Aniline	7.05	93	127277	23.34	ng	100
8) 2-Chlorophenol	7.18	128	79978	24.01	ng	99
9) Benzaldehyde	6.94	77	56141	25.81	ng	98
10) Phenol	7.00	94	99791	24.57	ng	99
11) bis(2-Chloroethyl)ether	7.11	93	81961	25.46	ng	99
12) 1,3-Dichlorobenzene	7.34	146	93547	24.57	ng	100
13) 1,4-Dichlorobenzene	7.42	146	94035	22.09	ng	100
14) 1,2-Dichlorobenzene	7.56	146	87123	23.59	ng	99
15) Benzyl Alcohol	7.52	79	67967	22.46	ng	100
16) 2,2'-oxybis(1-Chloropropan	7.66	45	116411	23.70	ng	99
17) 2-Methylphenol	7.62	107	70343	24.75	ng	97
18) Hexachloroethane	7.92	117	30454	22.92	ng	97
19) n-Nitroso-di-n-propylamine	7.78	70	65344	24.38	ng	99
20) 3+4-Methylphenols	7.77	107	88286	23.93	ng	98
22) Acetophenone	7.79	105	115295	24.41	ng	99
24) Nitrobenzene	7.96	77	78658	23.83	ng	99
25) Isophorone	8.20	82	170835	24.63	ng	99
26) 2-Nitrophenol	8.30	139	23771	18.81	ng	98
27) 2,4-Dimethylphenol	8.31	122	75473	24.31	ng	98
28) bis(2-Chloroethoxy)methane	8.41	93	98404	23.69	ng	99
29) 2,4-Dichlorophenol	8.52	162	62012	23.61	ng	99
30) 1,2,4-Trichlorobenzene	8.63	180	79500	23.58	ng	97
31) Naphthalene	8.71	128	264646	25.86	ng	99
32) Benzoic acid	8.35	122	20051m	21.35	ng	
33) 4-Chloroaniline	8.74	127	101985m	24.74	ng	
34) Hexachlorobutadiene	8.83	225	46484	25.17	ng	95
35) Caprolactam	9.07	113	18284	22.98	ng	# 65
36) 4-Chloro-3-methylphenol	9.21	107	73693	25.53	ng	100
37) 2-Methylnaphthalene	9.40	142	178451	24.65	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.58	216	80511	24.95	ng	99
40) Hexachlorocyclopentadiene	9.56	237	24646	16.58	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.68	196	48231	24.47	nq	99
43) 2,4,5-Trichlorophenol	9.71	196	54930	26.59	nq	97
45) 1,1'-Biphenyl	9.87	154	196869	25.58	nq	92
46) 2-Chloronaphthalene	9.90	162	167225	26.35	nq	98
47) 2-Nitroaniline	9.98	65	27696	20.07	nq	97
48) Acenaphthylene	10.32	152	278200	25.68	nq	99
49) Dimethylphthalate	10.15	163	186041	25.27	nq	99
50) 2,6-Dinitrotoluene	10.22	165	30052	23.40	nq	98
51) Acenaphthene	10.49	154	157063	25.03	nq	98
52) 3-Nitroaniline	10.39	138	37911	24.08	nq	95
53) 2,4-Dinitrophenol	10.49	184	2544	9.23	nq	# 1
54) Dibenzofuran	10.66	168	226400	25.57	nq	99
55) 4-Nitrophenol	10.52	139	25718	22.68	nq	95
56) 2,4-Dinitrotoluene	10.63	165	32774	20.10	nq	98
57) Fluorene	11.02	166	204581	26.01	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.78	232	39018	25.39	nq	97
59) Diethylphthalate	10.86	149	195870	27.06	nq	99
60) 4-Chlorophenyl-phenylether	10.99	204	87855	25.74	nq	96
61) 4-Nitroaniline	11.00	138	38566	24.71	nq	97
62) Azobenzene	11.15	77	174110	24.57	nq	97
64) 4,6-Dinitro-2-methylphenol	11.04	198	6682	12.64	nq	92
65) n-Nitrosodiphenylamine	11.11	169	159797	24.36	nq	99
66) 4-Bromophenyl-phenylether	11.51	248	52227	22.77	nq	# 88
67) Hexachlorobenzene	11.59	284	60900	25.50	nq	# 90
68) Atrazine	11.64	200	45351	24.04	nq	98
69) Pentachlorophenol	11.78	266	19764	22.24	nq	91
70) Phenanthrene	12.01	178	295245	24.94	nq	99
71) Anthracene	12.07	178	290879	25.43	nq	98
72) Carbazole	12.23	167	280893	26.01	nq	100
73) Di-n-butylphthalate	12.57	149	303600	27.13	nq	99
74) Fluoranthene	13.35	202	321367	25.82	nq	98
76) Benzidine	13.47	184	143126	23.69	nq	99
77) Pyrene	13.62	202	332670	25.86	nq	99
79) Butylbenzylphthalate	14.38	149	111243	25.35	nq	93
80) Benzo(a)anthracene	15.15	228	315351m	25.55	nq	
81) 3,3'-Dichlorobenzidine	15.10	252	102691m	26.32	nq	
82) Chrysene	15.20	228	298808m	25.81	nq	
83) Bis(2-ethylhexyl)phthalate	15.14	149	181677m	26.59	nq	
84) Di-n-octyl phthalate	16.03	149	281845m	26.22	nq	
85) Indeno(1,2,3-cd)pyrene	19.13	276	342116m	27.70	nq	
87) Benzo(b)fluoranthene	16.63	252	280136m	24.18	nq	
88) Benzo(k)fluoranthene	16.66	252	320078m	26.55	nq	
89) Benzo(a)pyrene	17.11	252	287142m	26.24	nq	
90) Dibenzo(a,h)anthracene	19.17	278	282058m	27.87	nq	
91) Benzo(g,h,i)perylene	19.71	276	296215m	27.68	nq	

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

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