

Data Path : Z:\HPCHEM1\BNA F\DATA\BF021316\
 Data File : BF085051.D
 Acq On : 12 Feb 2016 17:25
 Operator : SJ/IZ
 Sample : SSTDICC02.5
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC02.5

Manual Integrations
 APPROVED

Sohil
 2/15/2016 4:42:44 PM

Quant Time: Feb 12 23:28:46 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 12 23:09:50 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.87	152	142278	20.00	ng	-0.01
21) Naphthalene-d8	7.78	136	624458	20.00	ng	-0.01
38) Acenaphthene-d10	10.58	164	292421	20.00	ng	0.00
63) Phenanthrene-d10	12.98	188	550684	20.00	ng	-0.01
75) Chrysene-d12	17.17	240	383645	20.00	ng	-0.01
86) Perylene-d12	18.79	264	269375	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	4.41	112	43853m	4.88	ng	0.25
7) Phenol-d6	5.55	99	54366m	4.75	ng	0.11
23) Nitrobenzene-d5	6.73	82	53415	4.85	ng	0.01
41) 2,4,6-Tribromophenol	11.90	330	13695	4.49	ng	0.01
44) 2-Fluorobiphenyl	9.53	172	110934	5.58	ng	0.00
78) Terphenyl-d14	15.61	244	103986	6.04	ng	-0.01

Target Compounds

						Qvalue
8) 2-Chlorophenol	5.63	128	23660	2.29	ng	91
9) Benzaldehyde	5.35	77	15922	2.42	ng	# 77
10) Phenol	5.56	94	32645m	2.52	ng	
11) bis(2-Chloroethyl)ether	5.55	93	39821m	3.90	ng	
12) 1,3-Dichlorobenzene	5.82	146	23511	2.15	ng	97
13) 1,4-Dichlorobenzene	5.91	146	31302m	2.84	ng	
14) 1,2-Dichlorobenzene	6.12	146	28763	2.83	ng	100
15) Benzyl Alcohol	6.30	79	18343	2.37	ng	# 70
16) 2,2'-oxybis(1-Chloropropan	6.30	45	44194	2.63	ng	78
17) 2-Methylphenol	6.40	107	22682m	2.85	ng	
18) Hexachloroethane	6.58	117	10179	2.42	ng	# 85
19) n-Nitroso-di-n-propylamine	6.50	70	17421	2.43	ng	# 92
20) 3+4-Methylphenols	6.68	107	30643m	3.07	ng	
22) Acetophenone	6.52	105	40744	2.47	ng	# 98
24) Nitrobenzene	6.76	77	29397	2.45	ng	98
25) Isophorone	7.08	82	44482	2.09	ng	97
26) 2-Nitrophenol	7.26	139	12142	2.07	ng	95
27) 2,4-Dimethylphenol	7.42	122	21199	2.09	ng	92
28) bis(2-Chloroethoxy)methane	7.50	93	30608	2.40	ng	95
29) 2,4-Dichlorophenol	7.68	162	19001	2.20	ng	95
30) 1,2,4-Trichlorobenzene	7.71	180	24037	2.44	ng	# 96
31) Naphthalene	7.83	128	78799	2.50	ng	97
33) 4-Chloroaniline	8.02	127	28076	2.16	ng	# 92
34) Hexachlorobutadiene	8.01	225	12602	2.23	ng	95
36) 4-Chloro-3-methylphenol	8.86	107	21146	2.12	ng	89
37) 2-Methylnaphthalene	8.94	142	55193m	2.84	ng	
39) 1,2,4,5-Tetrachlorobenzene	9.20	216	23975	2.53	ng	98
43) 2,4,5-Trichlorophenol	9.55	196	15893m	2.59	ng	
45) 1,1'-Biphenyl	9.69	154	70017	2.67	ng	100
46) 2-Chloronaphthalene	9.70	162	49781	2.58	ng	98
47) 2-Nitroaniline	9.96	65	13122	2.12	ng	90
48) Acenaphthylene	10.35	152	77142	2.62	ng	96
49) Dimethylphthalate	10.22	163	56321	2.53	ng	# 96
50) 2,6-Dinitrotoluene	10.32	165	9993	2.08	ng	# 84

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51) Acenaphthene	10.64	154	47898	2.53	nq	98
52) 3-Nitroaniline	10.64	138	11054	2.02	nq	# 94
54) Dibenzofuran	10.92	168	64267	2.75	nq	95
56) 2,4-Dinitrotoluene	10.99	165	13574	2.17	nq	# 92
57) Fluorene	11.47	166	55039	2.81	nq	95
58) 2,3,4,6-Tetrachlorophenol	11.19	232	11686m	2.46	nq	
59) Diethylphthalate	11.36	149	56080	2.56	nq	99
60) 4-Chlorophenyl-phenylether	11.51	204	26004	2.76	nq	96
62) Azobenzene	11.76	77	53259	2.53	nq	96
65) n-Nitrosodiphenylamine	11.71	169	45814	2.65	nq	91
66) 4-Bromophenyl-phenylether	12.30	248	14512	2.40	nq	91
67) Hexachlorobenzene	12.35	284	15992	2.40	nq	91
68) Atrazine	12.63	200	12474	2.26	nq	90
70) Phenanthrene	13.03	178	81755	2.66	nq	95
71) Anthracene	13.11	178	84529	2.77	nq	99
72) Carbazole	13.43	167	64083	2.42	nq	99
73) Di-n-butylphthalate	14.03	149	95995	2.80	nq	97
74) Fluoranthene	14.96	202	78026	2.55	nq	98
77) Pyrene	15.31	202	78714	2.69	nq	99
79) Butylbenzylphthalate	16.45	149	30536	2.32	nq	92
80) Benzo(a)anthracene	17.15	228	56160	2.38	nq	98
82) Chrysene	17.20	228	57146	2.51	nq	99
83) Bis(2-ethylhexyl)phthalate	17.27	149	39453	2.37	nq	# 96
84) Di-n-octyl phthalate	18.02	149	55578	2.07	nq	# 90
85) Indeno(1,2,3-cd)pyrene	20.09	276	45757	2.53	nq	100
87) Benzo(b)fluoranthene	18.40	252	49805m	2.74	nq	
88) Benzo(k)fluoranthene	18.43	252	33674m	2.27	nq	
89) Benzo(a)pyrene	18.73	252	39987	2.61	nq	# 92
90) Dibenzo(a,h)anthracene	20.10	278	37269	2.82	nq	96
91) Benzo(g,h,i)perylene	20.47	276	39848	2.99	nq	99

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

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