

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031015\
 Data File : BF077679.D
 Acq On : 10 Mar 2015 23:51
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

mohammad
 3/12/2015 2:22:17 PM

Quant Time: Mar 11 02:46:19 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 11 00:54:19 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.18	152	47722	20.00	ng	0.00
21) Naphthalene-d8	8.76	136	200538	20.00	ng	0.00
38) Acenaphthene-d10	10.93	164	101461	20.00	ng	0.00
63) Phenanthrene-d10	12.77	188	227251	20.00	ng	0.00
75) Chrysene-d12	16.05	240	224954	20.00	ng	-0.03
86) Perylene-d12	17.75	264	221780	20.00	ng	-0.16

System Monitoring Compounds

5) 2-Fluorophenol	5.47	112	239544	83.80	ng	0.00
7) Phenol-d6	6.75	99	286301	77.90	ng	0.00
23) Nitrobenzene-d5	7.88	82	257945	79.99	ng	0.00
41) 2,4,6-Tribromophenol	11.91	330	42915	43.93	ng	0.00
44) 2-Fluorobiphenyl	10.11	172	558289	85.80	ng	0.00
78) Terphenyl-d14	14.77	244	660662	68.66	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.08	88	51579	42.52	ng	92
3) Pyridine	2.80	79	166294	47.92	ng	96
4) n-Nitrosodimethylamine	2.71	42	68096	45.13	ng	87
6) Aniline	6.77	93	200333	39.44	ng	# 73
8) 2-Chlorophenol	6.92	128	133374	39.55	ng	84
9) Benzaldehyde	6.63	77	104120	46.66	ng	93
10) Phenol	6.77	94	165756	38.01	ng	# 32
11) bis(2-Chloroethyl)ether	6.88	93	124048	39.59	ng	94
12) 1,3-Dichlorobenzene	7.11	146	143380	39.05	ng	97
13) 1,4-Dichlorobenzene	7.21	146	147716	39.54	ng	98
14) 1,2-Dichlorobenzene	7.39	146	141208	39.74	ng	99
15) Benzyl Alcohol	7.37	79	105768	37.86	ng	# 87
16) 2,2'-oxybis(1-Chloropropan	7.55	45	158726	35.43	ng	100
17) 2-Methylphenol	7.52	107	102632	38.94	ng	97
18) Hexachloroethane	7.81	117	47484	38.06	ng	# 78
19) n-Nitroso-di-n-propylamine	7.70	70	87791	37.61	ng	98
20) 3+4-Methylphenols	7.71	107	135609	39.06	ng	# 73
22) Acetophenone	7.70	105	183334	38.86	ng	# 82
24) Nitrobenzene	7.90	77	131501	38.76	ng	98
25) Isophorone	8.20	82	240983	37.67	ng	97
26) 2-Nitrophenol	8.29	139	55721	40.07	ng	98
27) 2,4-Dimethylphenol	8.36	122	120048	38.58	ng	98
28) bis(2-Chloroethoxy)methane	8.48	93	148223	36.67	ng	99
29) 2,4-Dichlorophenol	8.59	162	117004	40.06	ng	89
30) 1,2,4-Trichlorobenzene	8.69	180	118871	37.02	ng	98
31) Naphthalene	8.79	128	377500	37.64	ng	99
32) Benzoic acid	8.50	122	8942m	5.91	ng	
33) 4-Chloroaniline	8.86	127	173537	38.92	ng	# 94
34) Hexachlorobutadiene	8.94	225	66008	36.01	ng	99
35) Caprolactam	9.29	113	34411	38.60	ng	96
36) 4-Chloro-3-methylphenol	9.47	107	115076	39.19	ng	95
37) 2-Methylnaphthalene	9.65	142	258079	36.24	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.85	216	118304	40.64	ng	99
40) Hexachlorocyclopentadiene	9.84	237	30948	19.83	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	10.00	196	68047	35.30	nq	100
43) 2,4,5-Trichlorophenol	10.05	196	89912	43.61	nq	# 88
45) 1,1'-Biphenyl	10.23	154	311003	40.46	nq	97
46) 2-Chloronaphthalene	10.25	162	254946	42.29	nq	97
47) 2-Nitroaniline	10.38	65	68106	45.89	nq	97
48) Acenaphthylene	10.76	152	400719	41.99	nq	99
49) Dimethylphthalate	10.62	163	272884	38.26	nq	100
50) 2,6-Dinitrotoluene	10.69	165	62828	43.93	nq	97
51) Acenaphthene	10.97	154	230475	40.47	nq	99
52) 3-Nitroaniline	10.89	138	74866	45.40	nq	99
53) 2,4-Dinitrophenol	11.03	184	1094	2.51	nq	# 54
54) Dibenzofuran	11.19	168	352880	40.13	nq	99
55) 4-Nitrophenol	11.11	139	37418	28.21	nq	# 72
56) 2,4-Dinitrotoluene	11.18	165	81917	42.39	nq	# 88
57) Fluorene	11.61	166	289039	41.11	nq	99
58) 2,3,4,6-Tetrachlorophenol	11.34	232	29318	17.69	nq	98
59) Diethylphthalate	11.49	149	281140	39.98	nq	99
60) 4-Chlorophenyl-phenylether	11.62	204	133076	39.29	nq	92
61) 4-Nitroaniline	11.64	138	81318	51.45	nq	90
62) Azobenzene	11.82	77	258015	38.68	nq	96
64) 4,6-Dinitro-2-methylphenol	11.68	198	2981	6.75	nq	90
65) n-Nitrosodiphenylamine	11.77	169	249692	33.11	nq	99
66) 4-Bromophenyl-phenylether	12.23	248	89654	33.69	nq	94
67) Hexachlorobenzene	12.28	284	98095	34.34	nq	98
68) Atrazine	12.44	200	82221	33.54	nq	98
69) Pentachlorophenol	12.54	266	7939	5.29	nq	99
70) Phenanthrene	12.80	178	480100	36.45	nq	100
71) Anthracene	12.86	178	491037	37.57	nq	98
72) Carbazole	13.07	167	482198	38.80	nq	100
73) Di-n-butylphthalate	13.53	149	531139	36.39	nq	97
74) Fluoranthene	14.27	202	490848	34.12	nq	98
76) Benzidine	14.46	184	379715	49.96	nq	99
77) Pyrene	14.55	202	523493	38.04	nq	99
79) Butylbenzylphthalate	15.38	149	216314	38.21	nq	# 84
80) Benzo(a)anthracene	16.04	228	496182	37.63	nq	99
81) 3,3'-Dichlorobenzidine	16.02	252	180326	37.33	nq	# 97
82) Chrysene	16.08	228	470262	37.99	nq	98
83) Bis(2-ethylhexyl)phthalate	16.10	149	310573	34.21	nq	# 95
84) Di-n-octyl phthalate	16.88	149	529656	34.95	nq	96
85) Indeno(1,2,3-cd)pyrene	19.15	276	549406m	38.92	nq	
87) Benzo(b)fluoranthene	17.32	252	542121	42.04	nq	# 96
88) Benzo(k)fluoranthene	17.35	252	416313m	32.94	nq	
89) Benzo(a)pyrene	17.69	252	480073	39.08	nq	99
90) Dibenzo(a,h)anthracene	19.18	278	462534	39.48	nq	99
91) Benzo(g,h,i)perylene	19.56	276	465400	39.36	nq	96

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

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