

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
 Data File : BF085161.D
 Acq On : 3 Mar 2016 17:29
 Operator : SJ/UM
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

UMANGI
 3/4/2016 4:28:27 PM

Quant Time: Mar 04 00:24:40 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 04 00:11:57 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	188851	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	731936	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	347235	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	637488	20.00	ng	0.00
75) Chrysene-d12	14.15	240	392757	20.00	ng	0.00
86) Perylene-d12	15.61	264	312800	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.58	112	1208489	104.54	ng	0.00
7) Phenol-d6	6.61	99	1606173	106.08	ng	0.00
23) Nitrobenzene-d5	7.56	82	1659885	121.08	ng	0.01
41) 2,4,6-Tribromophenol	10.81	330	339112	99.96	ng	0.00
44) 2-Fluorobiphenyl	9.35	172	2464842	105.26	ng	0.00
78) Terphenyl-d14	13.09	244	2010188	119.45	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.66	88	327306	57.11	ng	100
3) Pyridine	3.44	79	805139	55.32	ng	100
4) n-Nitrosodimethylamine	3.40	42	363912	54.89	ng	96
6) Aniline	6.65	93	1188797	55.34	ng	96
8) 2-Chlorophenol	6.77	128	747215	55.71	ng	99
9) Benzaldehyde	6.53	77	354729	45.04	ng	99
10) Phenol	6.62	94	837342m	49.82	ng	
11) bis(2-Chloroethyl)ether	6.72	93	782487	58.23	ng	100
12) 1,3-Dichlorobenzene	6.93	146	828862	57.68	ng	98
13) 1,4-Dichlorobenzene	7.00	146	772301	52.88	ng	99
14) 1,2-Dichlorobenzene	7.16	146	737837	56.66	ng	99
15) Benzyl Alcohol	7.12	79	578622	52.18	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.26	45	1028315	52.07	ng	100
17) 2-Methylphenol	7.24	107	653823	58.37	ng	99
18) Hexachloroethane	7.50	117	311084	57.33	ng	99
19) n-Nitroso-di-n-propylamine	7.41	70	574443	57.97	ng	97
20) 3+4-Methylphenols	7.38	107	702012	52.56	ng	99
22) Acetophenone	7.40	105	946813	51.34	ng	97
24) Nitrobenzene	7.57	77	755639	52.68	ng	99
25) Isophorone	7.81	82	1508299	57.53	ng	99
26) 2-Nitrophenol	7.89	139	431591	67.25	ng	96
27) 2,4-Dimethylphenol	7.92	122	770901	62.20	ng	97
28) bis(2-Chloroethoxy)methane	8.01	93	789910	54.32	ng	99
29) 2,4-Dichlorophenol	8.13	162	605360	59.66	ng	98
30) 1,2,4-Trichlorobenzene	8.21	180	605259	54.50	ng	100
31) Naphthalene	8.29	128	1943700	52.78	ng	100
32) Benzoic acid	8.05	122	543400	88.93	ng	98
33) 4-Chloroaniline	8.33	127	815243	55.74	ng	100
34) Hexachlorobutadiene	8.41	225	338179	53.67	ng	99
35) Caprolactam	8.72	113	194562	60.74	ng	100
36) 4-Chloro-3-methylphenol	8.81	107	620543	55.19	ng	99
37) 2-Methylnaphthalene	8.98	142	1351643	59.43	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.14	216	522640	49.59	ng	99
40) Hexachlorocyclopentadiene	9.13	237	311637	52.51	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.26	196	427524	59.69	ng	99
43) 2,4,5-Trichlorophenol	9.29	196	402867	58.73	ng	98
45) 1,1'-Biphenyl	9.44	154	1609730	53.28	ng	100
46) 2-Chloronaphthalene	9.48	162	1291578	57.90	ng	98
47) 2-Nitroaniline	9.57	65	435544	62.46	ng	# 66
48) Acenaphthylene	9.89	152	1835096	54.58	ng	100
49) Dimethylphthalate	9.75	163	1504518	58.30	ng	99
50) 2,6-Dinitrotoluene	9.81	165	329738	61.58	ng	94
51) Acenaphthene	10.06	154	1130302	53.74	ng	99
52) 3-Nitroaniline	9.98	138	420214	66.34	ng	93
53) 2,4-Dinitrophenol	10.08	184	162760	92.56	ng	# 30
54) Dibenzofuran	10.23	168	1512649	55.25	ng	99
55) 4-Nitrophenol	10.13	139	330687	67.38	ng	85
56) 2,4-Dinitrotoluene	10.21	165	402921	55.95	ng	95
57) Fluorene	10.57	166	1207293	55.92	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.34	232	291457	56.90	ng	98
59) Diethylphthalate	10.45	149	1510780	61.48	ng	99
60) 4-Chlorophenyl-phenylether	10.56	204	568142	51.52	ng	99
61) 4-Nitroaniline	10.60	138	395942	67.24	ng	91
62) Azobenzene	10.72	77	1449184	58.20	ng	100
64) 4,6-Dinitro-2-methylphenol	10.62	198	250894	77.60	ng	88
65) n-Nitrosodiphenylamine	10.69	169	1163978	58.55	ng	99
66) 4-Bromophenyl-phenylether	11.05	248	351711	52.57	ng	96
67) Hexachlorobenzene	11.12	284	387830	52.14	ng	95
68) Atrazine	11.20	200	284597	47.50	ng	100
69) Pentachlorophenol	11.30	266	220744	51.20	ng	100
70) Phenanthrene	11.53	178	1735939	54.22	ng	100
71) Anthracene	11.59	178	2034207	57.38	ng	99
72) Carbazole	11.74	167	1746447	56.34	ng	99
73) Di-n-butylphthalate	12.07	149	2223097	60.47	ng	# 98
74) Fluoranthene	12.72	202	1853016	53.49	ng	99
76) Benzidine	12.84	184	712335	57.39	ng	97
77) Pyrene	12.95	202	1789454	63.98	ng	100
79) Butylbenzylphthalate	13.56	149	845379	71.66	ng	97
80) Benzo(a)anthracene	14.14	228	1405085	60.83	ng	99
81) 3,3'-Dichlorobenzidine	14.09	252	436584	60.32	ng	99
82) Chrysene	14.17	228	1318969	61.37	ng	99
83) Bis(2-ethylhexyl)phthalate	14.12	149	1041807	75.36	ng	100
84) Di-n-octyl phthalate	14.73	149	1696147	80.80	ng	99
85) Indeno(1,2,3-cd)pyrene	17.10	276	1150507	63.53	ng	100
87) Benzo(b)fluoranthene	15.18	252	1250038	64.32	ng	99
88) Benzo(k)fluoranthene	15.21	252	906440m	55.87	ng	
89) Benzo(a)pyrene	15.56	252	1032457	63.63	ng	98
90) Dibenzo(a,h)anthracene	17.11	278	958176	65.69	ng	99
91) Benzo(g,h,i)perylene	17.55	276	975411	65.69	ng	99

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

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