

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF083116\
 Data File : BF089934.D
 Acq On : 31 Aug 2016 5:01
 Operator : UM/SJ
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

SOHIL
 9/1/2016 11:51:59 AM

Quant Time: Aug 31 05:48:33 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF083116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 31 05:39:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.64	152	185003	20.00	ng	0.00
21) Naphthalene-d8	7.93	136	747986	20.00	ng	0.00
38) Acenaphthene-d10	9.67	164	372124	20.00	ng	0.00
63) Phenanthrene-d10	11.14	188	739036	20.00	ng	0.00
75) Chrysene-d12	13.76	240	560804	20.00	ng	0.00
86) Perylene-d12	15.10	264	460162	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.21	112	1233286	53.24	ng	0.00
7) Phenol-d6	6.28	99	1523318	53.65	ng	0.01
23) Nitrobenzene-d5	7.21	82	1411820	55.43	ng	0.00
41) 2,4,6-Tribromophenol	10.46	330	399735	56.55	ng	0.00
44) 2-Fluorobiphenyl	9.00	172	2121817	46.98	ng	0.00
78) Terphenyl-d14	12.73	244	2480047	54.31	ng	0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.10	88	314618	57.99	ng	99
3) Pyridine	2.73	79	925717	64.21	ng	96
4) n-Nitrosodimethylamine	2.70	42	435484	60.32	ng	100
6) Aniline	6.31	93	1063205	55.12	ng	# 73
8) 2-Chlorophenol	6.42	128	669697	52.81	ng	91
9) Benzaldehyde	6.18	77	521551	56.73	ng	95
10) Phenol	6.30	94	787972	47.73	ng	99
11) bis(2-Chloroethyl)ether	6.39	93	730871	58.90	ng	95
12) 1,3-Dichlorobenzene	6.58	146	819952	57.60	ng	98
13) 1,4-Dichlorobenzene	6.66	146	823562m	56.87	ng	
14) 1,2-Dichlorobenzene	6.81	146	662798	51.65	ng	98
15) Benzyl Alcohol	6.79	79	492498	54.56	ng	100
16) 2,2'-oxybis(1-Chloropropan	6.94	45	1356558	56.78	ng	98
17) 2-Methylphenol	6.91	107	583049	57.84	ng	98
18) Hexachloroethane	7.15	117	271952	54.90	ng	96
19) n-Nitroso-di-n-propylamine	7.08	70	499757	53.43	ng	# 93
20) 3+4-Methylphenols	7.07	107	715361	58.84	ng	# 70
22) Acetophenone	7.06	105	871682	52.73	ng	# 97
24) Nitrobenzene	7.23	77	820106	60.92	ng	91
25) Isophorone	7.48	82	1484700	56.91	ng	97
26) 2-Nitrophenol	7.54	139	388646	59.99	ng	95
27) 2,4-Dimethylphenol	7.60	122	725050	59.04	ng	98
28) bis(2-Chloroethoxy)methane	7.69	93	807459	51.03	ng	99
29) 2,4-Dichlorophenol	7.79	162	638679	58.32	ng	91
30) 1,2,4-Trichlorobenzene	7.87	180	691790	58.49	ng	99
31) Naphthalene	7.95	128	2111901	56.79	ng	99
32) Benzoic acid	7.76	122	565976	70.47	ng	97
33) 4-Chloroaniline	8.00	127	774366	51.91	ng	97
34) Hexachlorobutadiene	8.08	225	388577	56.06	ng	97
35) Caprolactam	8.39	113	207040	60.85	ng	93
36) 4-Chloro-3-methylphenol	8.49	107	619704	54.95	ng	84
37) 2-Methylnaphthalene	8.64	142	1370270	56.05	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.81	216	608307	55.34	ng	98
40) Hexachlorocyclopentadiene	8.80	237	370288	65.75	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.91	196	455832	60.45	ng	98
43) 2,4,5-Trichlorophenol	8.96	196	444115	59.61	ng #	86
45) 1,1'-Biphenyl	9.11	154	1762537	57.57	ng	95
46) 2-Chloronaphthalene	9.12	162	1140978	53.84	ng	98
47) 2-Nitroaniline	9.22	65	466624	68.10	ng #	85
48) Acenaphthylene	9.53	152	1834641	52.61	ng	100
49) Dimethylphthalate	9.42	163	1489896	57.55	ng	100
50) 2,6-Dinitrotoluene	9.47	165	402801	65.60	ng #	66
51) Acenaphthene	9.70	154	1244322	55.31	ng	100
52) 3-Nitroaniline	9.63	138	402574	62.02	ng	92
53) 2,4-Dinitrophenol	9.74	184	215243	86.49	ng #	74
54) Dibenzofuran	9.87	168	1505144	51.54	ng	97
55) 4-Nitrophenol	9.79	139	290240	55.75	ng #	60
56) 2,4-Dinitrotoluene	9.86	165	429215	56.01	ng #	82
57) Fluorene	10.22	166	1115429	49.90	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.00	232	388532	62.33	ng	94
59) Diethylphthalate	10.11	149	1519406	58.06	ng	96
60) 4-Chlorophenyl-phenylether	10.22	204	567658	53.34	ng #	90
61) 4-Nitroaniline	10.25	138	432816	66.72	ng	99
62) Azobenzene	10.38	77	1501632	59.35	ng	88
64) 4,6-Dinitro-2-methylphenol	10.27	198	304514	67.89	ng	81
65) n-Nitrosodiphenylamine	10.33	169	1213830	53.72	ng	99
66) 4-Bromophenyl-phenylether	10.70	248	389378	52.30	ng	94
67) Hexachlorobenzene	10.77	284	491515	56.76	ng #	71
68) Atrazine	10.87	200	411190	57.45	ng	97
69) Pentachlorophenol	10.95	266	281616	62.87	ng	98
70) Phenanthrene	11.17	178	1759851	49.00	ng	99
71) Anthracene	11.22	178	2175160	57.50	ng	99
72) Carbazole	11.37	167	1739990	50.32	ng	99
73) Di-n-butylphthalate	11.73	149	2387366	56.88	ng	99
74) Fluoranthene	12.34	202	2020845	53.03	ng	96
76) Benzidine	12.47	184	1132945	54.69	ng	97
77) Pyrene	12.57	202	2261753	57.90	ng	99
79) Butylbenzylphthalate	13.20	149	872097	56.45	ng	96
80) Benzo(a)anthracene	13.75	228	1938230	57.43	ng	99
81) 3,3'-Dichlorobenzidine	13.73	252	782891	64.12	ng #	95
82) Chrysene	13.78	228	1473639	49.15	ng	100
83) Bis(2-ethylhexyl)phthalate	13.77	149	1269051	57.85	ng	98
84) Di-n-octyl phthalate	14.38	149	2176319	62.48	ng	99
85) Indeno(1,2,3-cd)pyrene	16.33	276	1359168	59.15	ng	100
87) Benzo(b)fluoranthene	14.74	252	1702624m	59.60	ng	
88) Benzo(k)fluoranthene	14.77	252	1249187m	51.70	ng	
89) Benzo(a)pyrene	15.05	252	1418834	58.31	ng #	98
90) Dibenzo(a,h)anthracene	16.35	278	1110178	59.93	ng	97
91) Benzo(g,h,i)perylene	16.70	276	1132730	61.12	ng #	94

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

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