

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072016\
 Data File : BF089129.D
 Acq On : 20 Jul 2016 12:05
 Operator : UM/SJ
 Sample : SSTDIC02.5
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDIC02.5

Quant Time: Jul 20 16:43:25 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 20 15:28:22 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	47686	20.00	ng	0.00
21) Naphthalene-d8	7.88	136	211084	20.00	ng	0.00
38) Acenaphthene-d10	9.63	164	105204	20.00	ng	0.00
63) Phenanthrene-d10	11.11	188	209660	20.00	ng	0.00
75) Chrysene-d12	13.73	240	171307	20.00	ng	0.00
86) Perylene-d12	15.07	264	125025	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.21	112	15301	4.83	ng	0.02
7) Phenol-d6	6.26	99	22084	5.38	ng	-0.01
23) Nitrobenzene-d5	7.16	82	20297	5.04	ng	-0.01
41) 2,4,6-Tribromophenol	10.42	330	4505	4.69	ng	0.00
44) 2-Fluorobiphenyl	8.96	172	41846	5.99	ng	-0.01
78) Terphenyl-d14	12.69	244	37227	4.88	ng	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.06	88	3613	2.33	ng	92
3) Pyridine	2.73	79	7762	1.75	ng	# 69
4) n-Nitrosodimethylamine	2.66	42	2967	1.73	ng	# 88
6) Aniline	6.26	93	14582	2.48	ng	# 85
8) 2-Chlorophenol	6.39	128	8741	2.54	ng	85
9) Benzaldehyde	6.15	77	6351	2.34	ng	99
10) Phenol	6.28	94	12254	2.62	ng	93
11) bis(2-Chloroethyl)ether	6.34	93	9089	2.57	ng	99
12) 1,3-Dichlorobenzene	6.54	146	10369	2.74	ng	97
13) 1,4-Dichlorobenzene	6.62	146	10381	2.77	ng	99
14) 1,2-Dichlorobenzene	6.76	146	9817	2.75	ng	90
15) Benzyl Alcohol	6.76	79	7375	2.45	ng	100
16) 2,2'-oxybis(1-Chloropropan	6.89	45	10620	2.16	ng	82
17) 2-Methylphenol	6.89	107	7257	2.61	ng	99
18) Hexachloroethane	7.11	117	3800	2.64	ng	94
19) n-Nitroso-di-n-propylamine	7.02	70	6844	2.44	ng	# 87
20) 3+4-Methylphenols	7.05	107	9570	2.78	ng	99
22) Acetophenone	7.02	105	14589	2.79	ng	# 97
24) Nitrobenzene	7.19	77	10622	2.58	ng	95
25) Isophorone	7.43	82	18424	2.47	ng	99
26) 2-Nitrophenol	7.51	139	4533	2.65	ng	92
27) 2,4-Dimethylphenol	7.56	122	8649	2.54	ng	96
28) bis(2-Chloroethoxy)methane	7.64	93	11343	2.31	ng	96
29) 2,4-Dichlorophenol	7.76	162	7145	2.51	ng	97
30) 1,2,4-Trichlorobenzene	7.83	180	8217	2.65	ng	93
31) Naphthalene	7.91	128	30475	2.92	ng	100
33) 4-Chloroaniline	7.96	127	12016	2.58	ng	# 88
34) Hexachlorobutadiene	8.03	225	4557	2.72	ng	96
35) Caprolactam	8.30	113	2011	2.13	ng	# 84
36) 4-Chloro-3-methylphenol	8.47	107	8828	2.64	ng	89
37) 2-Methylnaphthalene	8.59	142	18225m	2.64	ng	
39) 1,2,4,5-Tetrachlorobenzene	8.76	216	10210	2.92	ng	95
42) 2,4,6-Trichlorophenol	8.89	196	5113	2.20	ng	97
43) 2,4,5-Trichlorophenol	8.92	196	5091	2.53	ng	# 80

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45) 1,1'-Biphenyl	9.06	154	24864	2.87	ng	96
46) 2-Chloronaphthalene	9.08	162	18436	2.71	ng #	91
47) 2-Nitroaniline	9.19	65	6062	2.55	ng #	80
48) Acenaphthylene	9.50	152	28251	2.60	ng	99
49) Dimethylphthalate	9.36	163	19891	2.58	ng #	96
50) 2,6-Dinitrotoluene	9.43	165	4557	2.66	ng	98
51) Acenaphthene	9.66	154	17078	2.56	ng	98
52) 3-Nitroaniline	9.60	138	5014	2.39	ng	98
54) Dibenzofuran	9.84	168	22984	2.63	ng	91
55) 4-Nitrophenol	9.79	139	2327m	1.50	ng	
56) 2,4-Dinitrotoluene	9.83	165	6416	2.90	ng #	88
57) Fluorene	10.17	166	21069	3.02	ng	98
58) 2,3,4,6-Tetrachlorophenol	9.96	232	3512	2.27	ng #	92
59) Diethylphthalate	10.06	149	20098	2.67	ng	98
60) 4-Chlorophenyl-phenylether	10.17	204	9353	2.93	ng #	75
61) 4-Nitroaniline	10.20	138	4720	2.29	ng	84
62) Azobenzene	10.33	77	22382	2.62	ng	91
65) n-Nitrosodiphenylamine	10.30	169	17133	2.46	ng	97
66) 4-Bromophenyl-phenylether	10.66	248	5090	2.41	ng #	81
67) Hexachlorobenzene	10.72	284	5961	2.59	ng #	78
68) Atrazine	10.82	200	5873	2.68	ng	95
70) Phenanthrene	11.13	178	28984	2.54	ng	98
71) Anthracene	11.18	178	32972	2.87	ng	99
72) Carbazole	11.34	167	26254	2.41	ng	98
73) Di-n-butylphthalate	11.68	149	34208	2.74	ng	99
74) Fluoranthene	12.31	202	31468	2.75	ng	93
76) Benzidine	12.43	184	14047	1.71	ng	98
77) Pyrene	12.53	202	33024	2.29	ng	99
79) Butylbenzylphthalate	13.17	149	13661	2.13	ng #	86
80) Benzo(a)anthracene	13.71	228	25278	2.31	ng	99
81) 3,3'-Dichlorobenzidine	13.68	252	7317	1.76	ng #	95
82) Chrysene	13.75	228	24530	2.25	ng	99
83) Bis(2-ethylhexyl)phthalate	13.73	149	18247	2.48	ng #	95
84) Di-n-octyl phthalate	14.33	149	29763	2.38	ng #	93
85) Indeno(1,2,3-cd)pyrene	16.29	276	16396	1.97	ng	99
87) Benzo(b)fluoranthene	14.71	252	23407m	2.84	ng	
88) Benzo(k)fluoranthene	14.73	252	16754m	2.55	ng	
89) Benzo(a)pyrene	15.02	252	17985	2.70	ng #	96
90) Dibenzo(a,h)anthracene	16.30	278	13274	2.40	ng #	90
91) Benzo(g,h,i)perylene	16.65	276	14014	2.46	ng	94

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

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