

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121418\
 Data File : BF111409.D
 Acq On : 14 Dec 2018 10:39
 Operator : JU/SJ
 Sample : SSTDICC025
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC025

Manual Integrations
 APPROVED

Sohil
 12/17/2018 2:36:06 PM

Quant Time: Dec 14 11:59:56 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 11:42:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	219127	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	936444	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	559302	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	999617	20.00	ng	0.00
76) Chrysene-d12	13.95	240	778770	20.00	ng	0.00
87) Perylene-d12	15.39	264	656246	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	704692	54.38	ng	0.00
7) Phenol-d6	6.43	99	905866	51.85	ng	0.00
23) Nitrobenzene-d5	7.36	82	847214	50.99	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	269492	54.66	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1905535	54.70	ng	0.00
79) Terphenyl-d14	12.90	244	1707003	48.24	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.51	88	165042	25.498	ng	98
3) Pyridine	3.24	79	416949	26.222	ng	95
4) n-Nitrosodimethylamine	3.17	42	191916	25.873	ng	96
6) Aniline	6.45	93	568228	26.884	ng	# 61
8) 2-Chlorophenol	6.58	128	409097	25.685	ng	97
9) Benzaldehyde	6.34	77	279379	25.695	ng	94
10) Phenol	6.45	94	506085	25.571	ng	78
11) bis(2-Chloroethyl)ether	6.53	93	387555	24.891	ng	99
12) 1,3-Dichlorobenzene	6.73	146	447955	25.980	ng	99
13) 1,4-Dichlorobenzene	6.81	146	453247	25.967	ng	97
14) 1,2-Dichlorobenzene	6.96	146	428219	26.519	ng	99
15) Benzyl Alcohol	6.93	79	346663	25.912	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.07	45	774845	25.688	ng	98
17) 2-Methylphenol	7.06	107	319516	24.827	ng	95
18) Hexachloroethane	7.31	117	166344	25.670	ng	93
19) n-Nitroso-di-n-propylamine	7.20	70	292590	25.202	ng	93
20) 3+4-Methylphenols	7.21	107	417970	27.233	ng	# 68
22) Acetophenone	7.20	105	564327	25.943	ng	# 91
24) Nitrobenzene	7.37	77	421199	24.612	ng	92
25) Isophorone	7.61	82	690981	24.484	ng	96
26) 2-Nitrophenol	7.69	139	222565	26.458	ng	94
27) 2,4-Dimethylphenol	7.73	122	317065	24.830	ng	96
28) bis(2-Chloroethoxy)methane	7.83	93	481605	24.862	ng	100
29) 2,4-Dichlorophenol	7.94	162	340312	25.549	ng	95
30) 1,2,4-Trichlorobenzene	8.02	180	357561	26.125	ng	99
31) Naphthalene	8.10	128	1153509	26.439	ng	96
32) Benzoic acid	7.84	122	272234m	25.646	ng	
33) 4-Chloroaniline	8.15	127	508133	27.271	ng	97
34) Hexachlorobutadiene	8.22	225	198039	26.083	ng	97
35) Caprolactam	8.51	113	124997	26.781	ng	98
36) 4-Chloro-3-methylphenol	8.63	107	373035	26.280	ng	99
37) 2-Methylnaphthalene	8.79	142	932689	33.070	ng	97
38) 1-Methylnaphthalene	8.89	142	894135	32.848	ng	98
40) 1,2,4,5-Tetrachlorobenzene	8.96	216	406911	25.883	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.95	237	207246	25.245	nq	99
43) 2,4,6-Trichlorophenol	9.07	196	286274	25.441	nq	98
44) 2,4,5-Trichlorophenol	9.11	196	297694	25.032	nq	98
46) 1,1'-Biphenyl	9.25	154	1248119	26.155	nq	96
47) 2-Chloronaphthalene	9.27	162	951879	26.046	nq	97
48) 2-Nitroaniline	9.37	65	301886	25.420	nq	95
49) Acenaphthylene	9.69	152	1421721	26.627	nq	95
50) Dimethylphthalate	9.55	163	1045437	25.993	nq	99
51) 2,6-Dinitrotoluene	9.61	165	241642	27.104	nq	92
52) Acenaphthene	9.86	154	885824	26.303	nq	98
53) 3-Nitroaniline	9.78	138	290891	27.146	nq	97
54) 2,4-Dinitrophenol	9.89	184	87113	24.906	nq	# 84
55) Dibenzofuran	10.03	168	1306162	26.866	nq	96
56) 4-Nitrophenol	9.95	139	209190	27.108	nq	91
57) 2,4-Dinitrotoluene	10.02	165	320910	28.086	nq	90
58) Fluorene	10.38	166	930344	26.750	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.16	232	240998	26.233	nq	97
60) Diethylphthalate	10.25	149	1088914	26.893	nq	98
61) 4-Chlorophenyl-phenylether	10.37	204	460322	26.698	nq	98
62) 4-Nitroaniline	10.39	138	286202	27.786	nq	95
63) Azobenzene	10.53	77	1066168	26.115	nq	95
65) 4,6-Dinitro-2-methylphenol	10.42	198	144232	25.133	nq	94
66) n-Nitrosodiphenylamine	10.49	169	899408	25.294	nq	98
67) 4-Bromophenyl-phenylether	10.86	248	281235	25.272	nq	93
68) Hexachlorobenzene	10.93	284	293632	25.625	nq	94
69) Atrazine	11.02	200	276649	26.900	nq	98
70) Pentachlorophenol	11.12	266	175099	22.962	nq	95
71) Phenanthrene	11.34	178	1351341	26.154	nq	95
72) Anthracene	11.39	178	1387739	26.378	nq	95
73) Carbazole	11.54	167	1444202	26.549	nq	95
74) Di-n-butylphthalate	11.88	149	1641233	26.498	nq	# 96
75) Fluoranthene	12.53	202	1385557	27.470	nq	97
77) Benzidine	12.64	184	717869	48.846	nq	98
78) Pyrene	12.75	202	1433595	23.852	nq	96
80) Butylbenzylphthalate	13.37	149	695130	24.088	nq	92
81) Benzo(a)anthracene	13.94	228	1087042	24.947	nq	99
82) 3,3'-Dichlorobenzidine	13.90	252	468484	28.429	nq	97
83) Chrysene	13.98	228	1162986	26.161	nq	98
84) Bis(2-ethylhexyl)phthalate	13.94	149	875153	24.807	nq	99
85) Di-n-octyl phthalate	14.54	149	1519534	26.232	nq	97
86) Indeno(1,2,3-cd)pyrene	16.80	276	737144	21.201	nq	98
88) Benzo(b)fluoranthene	14.97	252	982122	26.319	nq	98
89) Benzo(k)fluoranthene	15.00	252	990005	26.579	nq	98
90) Benzo(a)pyrene	15.33	252	895830	25.783	nq	99
91) Dibenzo(a,h)anthracene	16.82	278	630601	21.814	nq	96
92) Benzo(g,h,i)perylene	17.22	276	583462	20.188	nq	95

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

